



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 20, 2016 – 01:16 PM GMT

PDB ID : 4WT1
Title : Complex of 70S ribosome with tRNA-Phe and mRNA with A-A mismatch in the second position in the A-site
Authors : Rozov, A.; Demeshkina, N.; Yusupov, M.; Yusupova, G.
Deposited on : 2014-10-29
Resolution : 3.05 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026982
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20026982

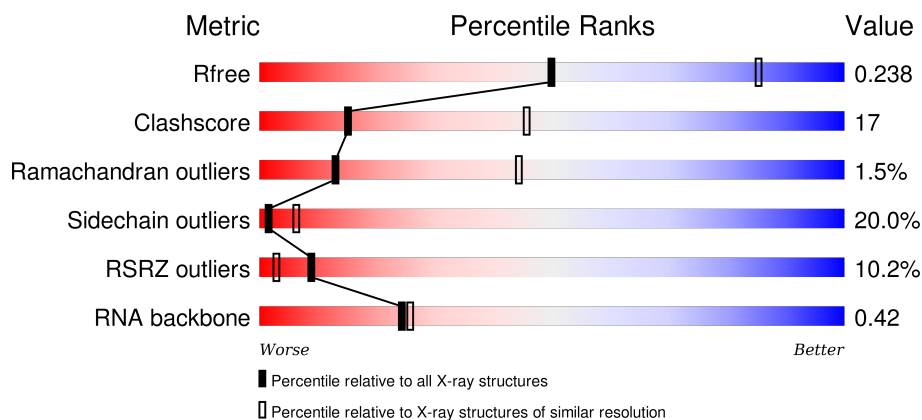
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.05 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	91344	1191 (3.10-3.02)
Clashscore	102246	1303 (3.10-3.02)
Ramachandran outliers	100387	1254 (3.10-3.02)
Sidechain outliers	100360	1254 (3.10-3.02)
RSRZ outliers	91569	1197 (3.10-3.02)
RNA backbone	2183	1035 (3.50-2.62)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	13	1522	<div> <div>3%</div> <div> <div></div> <div>33%</div> <div>43%</div> <div>19%</div> <div>••</div> </div> </div>
1	1G	1522	<div> <div>5%</div> <div> <div></div> <div>33%</div> <div>46%</div> <div>17%</div> <div>••</div> </div> </div>
2	12	256	<div> <div>7%</div> <div> <div></div> <div>45%</div> <div>38%</div> <div>10%</div> <div>7%</div> </div> </div>
2	1E	256	<div> <div>3%</div> <div> <div></div> <div>41%</div> <div>39%</div> <div>12%</div> <div>7%</div> </div> </div>

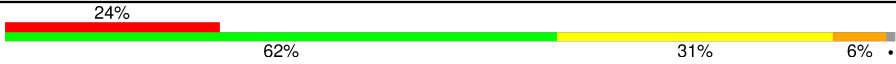

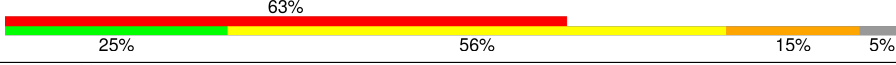

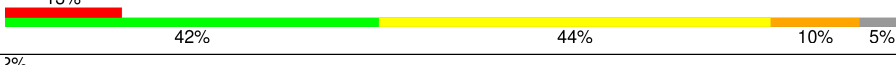
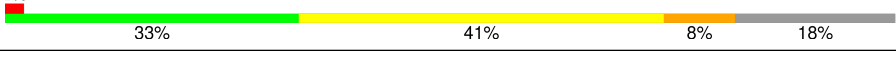
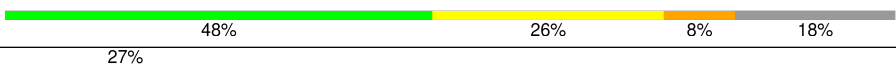
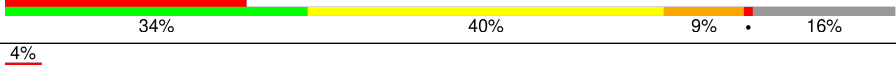
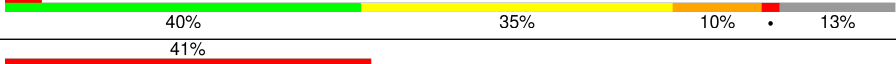



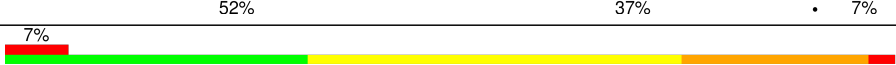
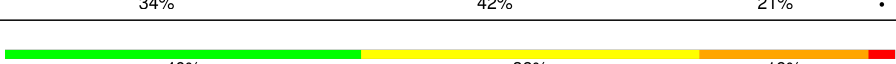
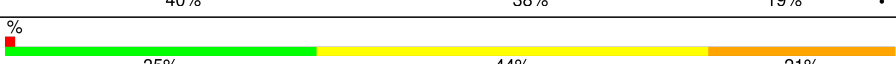
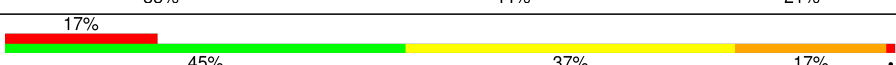
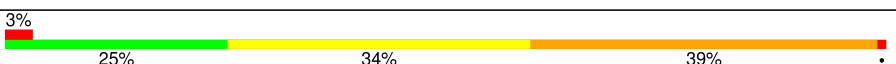
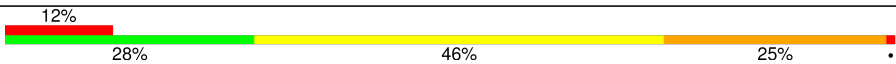
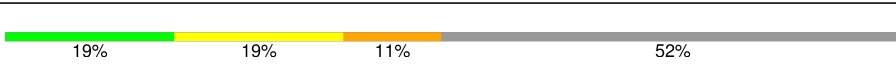

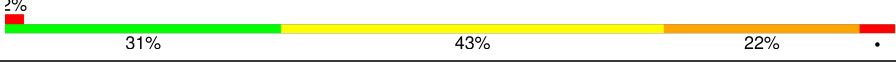
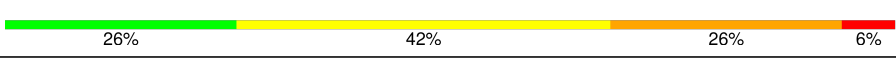

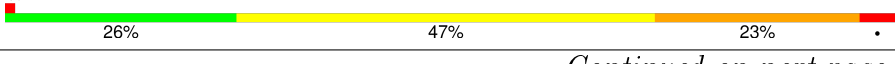

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Mol	Chain	Length	Quality of chain
3	22	239	
3	2E	239	
4	32	209	
4	3E	209	
5	42	162	
5	4E	162	
6	52	101	
6	5E	101	
7	62	156	
7	6E	156	
8	72	138	
8	7E	138	
9	82	128	
9	8E	128	
10	1A	105	
10	1I	105	
11	2A	129	
11	2I	129	
12	3A	132	
12	3I	132	
13	4A	126	
13	4I	126	
14	5A	61	
14	5I	61	
15	6A	89	

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Mol	Chain	Length	Quality of chain
15	6I	89	
16	7A	88	
16	7I	88	
17	8A	105	
17	8I	105	
18	9A	88	
18	9I	88	
19	AA	93	
19	AI	93	
20	BA	106	
20	BI	106	
21	1B	27	
21	1F	27	
22	1K	76	
23	2K	77	
23	2L	77	
24	1L	76	
24	3K	76	
24	3L	76	
25	4K	27	
25	4L	27	
26	14	2917	
26	1H	2917	
27	16	122	
27	1J	122	

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Mol	Chain	Length	Quality of chain
28	11	276	
28	19	276	
29	21	206	
29	29	206	
30	31	210	
30	39	210	
31	41	182	
31	49	182	
32	51	180	
32	59	180	
33	61	148	
33	69	148	
34	15	140	
34	58	140	
35	25	122	
35	68	122	
36	35	150	
36	78	150	
37	45	141	
37	88	141	
38	55	118	
38	98	118	
39	65	112	
39	A8	112	
40	75	146	

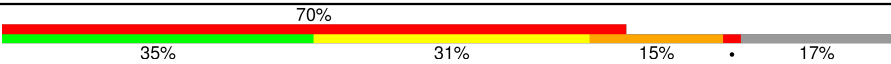


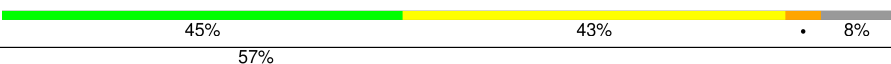

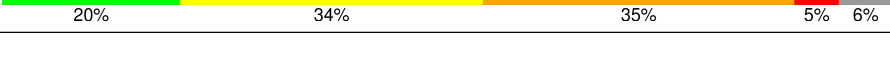
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Mol	Chain	Length	Quality of chain
40	B8	146	
41	85	118	
41	C8	118	
42	95	101	
42	D8	101	
43	A5	113	
43	E8	113	
44	B5	96	
44	F8	96	
45	C5	110	
45	G8	110	
46	D5	206	
46	H8	206	
47	E5	85	
47	I8	85	
48	F5	98	
48	J8	98	
49	G5	72	
49	K8	72	
50	H5	60	
50	L8	60	
51	I5	71	
51	M8	71	
52	J5	60	
52	N8	60	

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Mol	Chain	Length	Quality of chain
53	K5	54	
53	O8	54	
54	L5	49	
54	P8	49	
55	M5	65	
55	Q8	65	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
56	MG	13	1602	-	-	-	X
56	MG	13	1621	-	-	-	X
56	MG	13	1641	-	-	-	X
56	MG	13	1652	-	-	-	X
56	MG	13	1664	-	-	-	X
56	MG	13	1666	-	-	-	X
56	MG	13	1680	-	-	-	X
56	MG	13	1686	-	-	-	X
56	MG	13	1689	-	-	-	X
56	MG	13	1706	-	-	-	X
56	MG	13	1721	-	-	-	X
56	MG	13	1752	-	-	-	X
56	MG	14	3007	-	-	-	X
56	MG	14	3055	-	-	-	X
56	MG	14	3060	-	-	-	X
56	MG	14	3091	-	-	-	X
56	MG	14	3146	-	-	-	X
56	MG	14	3169	-	-	-	X
56	MG	14	3170	-	-	-	X
56	MG	14	3194	-	-	-	X
56	MG	14	3195	-	-	-	X
56	MG	14	3199	-	-	-	X
56	MG	14	3202	-	-	-	X
56	MG	14	3205	-	-	-	X
56	MG	14	3209	-	-	-	X
56	MG	14	3211	-	-	-	X
56	MG	14	3216	-	-	-	X

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
56	MG	14	3217	-	-	-	X
56	MG	14	3227	-	-	-	X
56	MG	14	3233	-	-	-	X
56	MG	14	3248	-	-	-	X
56	MG	14	3253	-	-	-	X
56	MG	14	3254	-	-	-	X
56	MG	14	3264	-	-	-	X
56	MG	14	3266	-	-	-	X
56	MG	14	3272	-	-	-	X
56	MG	14	3274	-	-	-	X
56	MG	14	3306	-	-	-	X
56	MG	14	3308	-	-	-	X
56	MG	14	3382	-	-	-	X
56	MG	16	206	-	-	-	X
56	MG	16	207	-	-	-	X
56	MG	16	209	-	-	-	X
56	MG	1G	1614	-	-	-	X
56	MG	1G	1627	-	-	-	X
56	MG	1G	1645	-	-	-	X
56	MG	1G	1659	-	-	-	X
56	MG	1G	1660	-	-	-	X
56	MG	1G	1664	-	-	-	X
56	MG	1H	3004	-	-	-	X
56	MG	1H	3011	-	-	-	X
56	MG	1H	3013	-	-	-	X
56	MG	1H	3025	-	-	-	X
56	MG	1H	3037	-	-	-	X
56	MG	1H	3052	-	-	-	X
56	MG	1H	3056	-	-	-	X
56	MG	1H	3063	-	-	-	X
56	MG	1H	3069	-	-	-	X
56	MG	1H	3081	-	-	-	X
56	MG	1H	3084	-	-	-	X
56	MG	1H	3089	-	-	-	X
56	MG	1H	3110	-	-	-	X
56	MG	1H	3116	-	-	-	X
56	MG	1H	3139	-	-	-	X
56	MG	1H	3160	-	-	-	X
56	MG	1H	3170	-	-	-	X
56	MG	1H	3173	-	-	-	X
56	MG	1H	3178	-	-	-	X
56	MG	1H	3186	-	-	-	X

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
56	MG	1H	3219	-	-	-	X
56	MG	1H	3222	-	-	-	X
56	MG	1H	3234	-	-	-	X
56	MG	1H	3235	-	-	-	X
56	MG	1H	3240	-	-	-	X
56	MG	1H	3245	-	-	-	X
56	MG	1H	3254	-	-	-	X
56	MG	1H	3255	-	-	-	X
56	MG	1H	3258	-	-	-	X
56	MG	1H	3260	-	-	-	X
56	MG	1H	3277	-	-	-	X
56	MG	1H	3298	-	-	-	X
56	MG	1H	3304	-	-	-	X
56	MG	1H	3332	-	-	-	X
56	MG	1J	207	-	-	-	X
56	MG	29	301	-	-	-	X
56	MG	2K	105	-	-	-	X
56	MG	39	303	-	-	-	X
56	MG	L8	101	-	-	-	X
57	ZN	32	302	-	-	-	X
57	ZN	3E	301	-	-	-	X

2 Entry composition

There are 58 unique types of molecules in this entry. The entry contains 299951 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a RNA chain called 16S ribosomal RNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	13	1498	Total	C	N	O	P	0	0	0
			32207	14334	5973	10402	1498			
1	1G	1497	Total	C	N	O	P	0	0	0
			32182	14324	5968	10394	1496			

- Molecule 2 is a protein called 30S ribosomal protein S2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	1E	237	Total	C	N	O	S	0	0	0
			1924	1228	344	347	5			
2	12	237	Total	C	N	O	S	0	0	0
			1924	1228	344	347	5			

- Molecule 3 is a protein called 30S ribosomal protein S3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	2E	205	Total	C	N	O	S	0	0	0
			1605	1011	313	280	1			
3	22	206	Total	C	N	O	S	0	0	0
			1612	1016	314	281	1			

- Molecule 4 is a protein called 30S ribosomal protein S4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	3E	208	Total	C	N	O	S	0	0	0
			1702	1066	339	290	7			
4	32	208	Total	C	N	O	S	0	0	0
			1702	1066	339	290	7			

- Molecule 5 is a protein called 30S ribosomal protein S5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	4E	151	Total	C	N	O	S	0	0	0
			1155	729	218	204	4			
5	42	151	Total	C	N	O	S	0	0	0
			1155	729	218	204	4			

- Molecule 6 is a protein called 30S ribosomal protein S6.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
6	5E	101	Total	C	N	O	S	0	0	0
			842	531	155	153	3			
6	52	101	Total	C	N	O	S	0	0	0
			842	531	155	153	3			

- Molecule 7 is a protein called 30S ribosomal protein S7.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
7	6E	155	Total	C	N	O	S	0	0	0
			1256	781	252	217	6			
7	62	155	Total	C	N	O	S	0	0	0
			1256	781	252	217	6			

- Molecule 8 is a protein called 30S ribosomal protein S8.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
8	7E	138	Total	C	N	O	S	0	0	0
			1115	705	215	192	3			
8	72	138	Total	C	N	O	S	0	0	0
			1115	705	215	192	3			

- Molecule 9 is a protein called 30S ribosomal protein S9.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
9	8E	127	Total	C	N	O		0	0	0
			1009	639	197	173				
9	82	127	Total	C	N	O		0	0	0
			1009	639	197	173				

- Molecule 10 is a protein called 30S ribosomal protein S10.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
10	1I	99	Total	C	N	O	S	0	0	0
			801	504	157	139	1			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
10	1A	99	Total	C	N	O	S	0	0	0
			801	504	157	139	1			

- Molecule 11 is a protein called 30S ribosomal protein S11.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
11	2I	116	Total	C	N	O	S	0	0	0
			864	537	164	160	3			
11	2A	116	Total	C	N	O	S	0	0	0
			864	537	164	160	3			

- Molecule 12 is a protein called 30S ribosomal protein S12.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
12	3I	125	Total	C	N	O	S	0	0	0
			975	614	196	164	1			
12	3A	125	Total	C	N	O	S	0	0	0
			975	614	196	164	1			

- Molecule 13 is a protein called 30S ribosomal protein S13.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
13	4I	118	Total	C	N	O	S	0	0	0
			938	580	193	163	2			
13	4A	117	Total	C	N	O	S	0	0	0
			933	577	192	162	2			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
4I	119	ALA	GLY	conflict	UNP P80377
4A	119	ALA	GLY	conflict	UNP P80377

- Molecule 14 is a protein called 30S ribosomal protein S14 type Z.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
14	5I	60	Total	C	N	O	S	0	0	0
			491	312	104	71	4			
14	5A	58	Total	C	N	O	S	0	0	0
			475	303	99	69	4			

- Molecule 15 is a protein called 30S ribosomal protein S15.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
15	6I	88	Total	C	N	O	S	0	0	0
			733	459	147	125	2			
15	6A	88	Total	C	N	O	S	0	0	0
			733	459	147	125	2			

- Molecule 16 is a protein called 30S ribosomal protein S16.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
16	7I	84	Total	C	N	O	S	0	0	0
			705	446	140	118	1			
16	7A	84	Total	C	N	O	S	0	0	0
			705	446	140	118	1			

- Molecule 17 is a protein called 30S ribosomal protein S17.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
17	8I	100	Total	C	N	O	S	0	0	0
			834	534	155	143	2			
17	8A	100	Total	C	N	O	S	0	0	0
			834	534	155	143	2			

- Molecule 18 is a protein called 30S ribosomal protein S18.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
18	9I	72	Total	C	N	O	0	0	0
			590	376	117	97			
18	9A	72	Total	C	N	O	0	0	0
			590	376	117	97			

- Molecule 19 is a protein called 30S ribosomal protein S19.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
19	AI	81	Total	C	N	O	S	0	0	0
			647	413	119	113	2			
19	AA	78	Total	C	N	O	S	0	0	0
			624	398	115	109	2			

- Molecule 20 is a protein called 30S ribosomal protein S20.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
20	BI	99	Total	C	N	O	S	0	0	0
			762	470	162	128	2			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
20	BA	99	Total	C	N	O	S	0	0	0
			762	470	162	128	2			

- Molecule 21 is a protein called 30S ribosomal protein Thx.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
21	1F	25	Total	C	N	O	0	0	0
			217	134	52	31			
21	1B	25	Total	C	N	O	0	0	0
			217	134	52	31			

- Molecule 22 is a RNA chain called tRNA-Phe.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
22	1K	76	Total	C	N	O	P	S	0	0	0
			1628	731	290	530	75	2			

- Molecule 23 is a RNA chain called tRNA-fMet.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
23	2K	77	Total	C	N	O	P	S	0	0	0
			1646	735	298	535	77	1			
23	2L	77	Total	C	N	O	P	S	0	0	0
			1646	735	298	535	77	1			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
2K	18	C	U	conflict	GB 675818144
2L	18	C	U	conflict	GB 675818144

- Molecule 24 is a RNA chain called tRNA-Phe.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
24	3K	76	Total	C	N	O	P	0	0	0
			1619	723	290	531	75			
24	1L	76	Total	C	N	O	P	0	0	0
			1619	723	290	531	75			
24	3L	76	Total	C	N	O	P	0	0	0
			1619	723	290	531	75			

- Molecule 25 is a RNA chain called mRNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
25	4K	13	Total	C	N	O	P	0	0	0
			281	127	57	84	13			
25	4L	9	Total	C	N	O	P	0	0	0
			193	87	37	60	9			

- Molecule 26 is a RNA chain called 23S ribosomal RNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
26	1H	2912	Total	C	N	O	P	0	0	0
			62707	27911	11722	20163	2911			
26	14	2909	Total	C	N	O	P	0	0	0
			62647	27884	11716	20139	2908			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
1H	161	U	-	insertion	GB 48268
1H	493	G	-	insertion	GB 48268
1H	1228	G	-	insertion	GB 48268
14	161	U	-	insertion	GB 48268
14	493	G	-	insertion	GB 48268
14	1228	G	-	insertion	GB 48268

- Molecule 27 is a RNA chain called 5S ribosomal RNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
27	16	122	Total	C	N	O	P	0	0	0
			2617	1166	486	844	121			
27	1J	122	Total	C	N	O	P	0	0	0
			2617	1166	486	844	121			

- Molecule 28 is a protein called 50S ribosomal protein L2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
28	11	272	Total	C	N	O	S	0	0	0
			2115	1335	420	357	3			
28	19	273	Total	C	N	O	S	0	0	0
			2120	1338	421	358	3			

- Molecule 29 is a protein called 50S ribosomal protein L3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
29	21	205	Total	C	N	O	S	0	0	0
			1568	991	300	271	6			
29	29	205	Total	C	N	O	S	0	0	0
			1568	991	300	271	6			

- Molecule 30 is a protein called 50S ribosomal protein L4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
30	31	202	Total	C	N	O	S	0	0	0
			1585	1011	297	275	2			
30	39	208	Total	C	N	O	S	0	0	0
			1627	1037	304	283	3			

- Molecule 31 is a protein called 50S ribosomal protein L5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
31	41	181	Total	C	N	O	S	0	0	0
			1473	942	268	259	4			
31	49	181	Total	C	N	O	S	0	0	0
			1473	942	268	259	4			

- Molecule 32 is a protein called 50S ribosomal protein L6.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
32	51	174	Total	C	N	O	S	0	0	0
			1336	848	251	236	1			
32	59	169	Total	C	N	O	S	0	0	0
			1299	823	244	231	1			

- Molecule 33 is a protein called 50S ribosomal protein L9.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
33	61	146	Total	C	N	O	S	0	0	0
			1136	726	201	208	1			
33	69	146	Total	C	N	O	S	0	0	0
			1136	726	201	208	1			

- Molecule 34 is a protein called 50S ribosomal protein L13.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
34	58	138	Total	C	N	O	S	0	0	0
			1104	712	206	182	4			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
34	15	138	Total	C	N	O	S	0	0	0
			1104	712	206	182	4			

- Molecule 35 is a protein called 50S ribosomal protein L14.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
35	68	122	Total	C	N	O	S	0	0	0
			932	588	171	169	4			
35	25	122	Total	C	N	O	S	0	0	0
			932	588	171	169	4			

- Molecule 36 is a protein called 50S ribosomal protein L15.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
36	78	150	Total	C	N	O	S	0	0	0
			1144	712	232	197	3			
36	35	150	Total	C	N	O	S	0	0	0
			1144	712	232	197	3			

- Molecule 37 is a protein called 50S ribosomal protein L16.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
37	88	138	Total	C	N	O	S	0	0	0
			1086	693	208	179	6			
37	45	141	Total	C	N	O	S	0	0	0
			1121	715	212	187	7			

- Molecule 38 is a protein called 50S ribosomal protein L17.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
38	98	118	Total	C	N	O	S	0	0	0
			967	604	203	159	1			
38	55	117	Total	C	N	O	S	0	0	0
			959	599	202	158				

- Molecule 39 is a protein called 50S ribosomal protein L18.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
39	A8	111	Total	C	N	O	0	0	0
			881	556	176	149			
39	65	111	Total	C	N	O	0	0	0
			881	556	176	149			

- Molecule 40 is a protein called 50S ribosomal protein L19.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
40	B8	137	Total	C	N	O	S	0	0	0
			1141	710	234	196	1			
40	75	137	Total	C	N	O	S	0	0	0
			1141	710	234	196	1			

- Molecule 41 is a protein called 50S ribosomal protein L20.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
41	C8	117	Total	C	N	O	S	0	0	0
			963	610	202	150	1			
41	85	117	Total	C	N	O	S	0	0	0
			963	610	202	150	1			

- Molecule 42 is a protein called 50S ribosomal protein L21.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
42	D8	101	Total	C	N	O	S	0	0	0
			778	501	142	134	1			
42	95	101	Total	C	N	O	S	0	0	0
			778	501	142	134	1			

- Molecule 43 is a protein called 50S ribosomal protein L22.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
43	E8	113	Total	C	N	O	S	0	0	0
			899	566	177	154	2			
43	A5	113	Total	C	N	O	S	0	0	0
			899	566	177	154	2			

- Molecule 44 is a protein called 50S ribosomal protein L23.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
44	F8	94	Total	C	N	O	S	0	0	0
			742	482	134	125	1			
44	B5	92	Total	C	N	O		0	0	0
			725	471	131	123				

- Molecule 45 is a protein called 50S ribosomal protein L24.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
45	G8	104	Total	C	N	O	S	0	0	0
			791	510	149	127	5			
45	C5	104	Total	C	N	O	S	0	0	0
			794	510	152	127	5			

- Molecule 46 is a protein called 50S ribosomal protein L25.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
46	H8	175	Total	C	N	O	S	0	0	0
			1397	892	251	251	3			
46	D5	179	Total	C	N	O	S	0	0	0
			1428	911	255	259	3			

- Molecule 47 is a protein called 50S ribosomal protein L27.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
47	I8	80	Total	C	N	O	S	0	0	0
			626	388	132	105	1			
47	E5	76	Total	C	N	O	S	0	0	0
			606	376	128	101	1			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
I8	6	ALA	GLY	conflict	UNP P60493
I8	8	ALA	GLY	conflict	UNP P60493
E5	6	ALA	GLY	conflict	UNP P60493
E5	8	ALA	GLY	conflict	UNP P60493

- Molecule 48 is a protein called 50S ribosomal protein L28.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
48	J8	97	Total	C	N	O	S	0	0	0
			762	481	150	130	1			
48	F5	97	Total	C	N	O	S	0	0	0
			762	481	150	130	1			

- Molecule 49 is a protein called 50S ribosomal protein L29.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
49	K8	67	Total	C	N	O	S	0	0	0
			563	349	114	99	1			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
49	G5	66	Total	C	N	O	S	0	0	0
			558	346	113	98	1			

- Molecule 50 is a protein called 50S ribosomal protein L30.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
50	L8	57	Total	C	N	O		0	0	0
			452	288	88	76				
50	H5	59	Total	C	N	O		0	0	0
			468	298	90	80				

- Molecule 51 is a protein called 50S ribosomal protein L31.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
51	M8	66	Total	C	N	O	S	0	0	0
			533	335	96	97	5			
51	I5	63	Total	C	N	O	S	0	0	0
			515	326	93	91	5			

- Molecule 52 is a protein called 50S ribosomal protein L32.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
52	N8	58	Total	C	N	O	S	0	0	0
			453	285	89	74	5			
52	J5	56	Total	C	N	O	S	0	0	0
			434	272	87	70	5			

- Molecule 53 is a protein called 50S ribosomal protein L33.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
53	O8	45	Total	C	N	O	S	0	0	0
			389	241	79	65	4			
53	K5	45	Total	C	N	O	S	0	0	0
			389	241	79	65	4			

- Molecule 54 is a protein called 50S ribosomal protein L34.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
54	P8	45	Total	C	N	O	S	0	0	0
			391	240	97	52	2			
54	L5	46	Total	C	N	O	S	0	0	0
			398	245	98	53	2			

- Molecule 55 is a protein called 50S ribosomal protein L35.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
55	Q8	61	Total	C	N	O	S	0	0	0
			485	309	99	75	2			
55	M5	60	Total	C	N	O	S	0	0	0
			477	303	98	74	2			

- Molecule 56 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
56	45	2	Total	Mg	0	0
			2	2		
56	P8	1	Total	Mg	0	0
			1	1		
56	85	1	Total	Mg	0	0
			1	1		
56	32	1	Total	Mg	0	0
			1	1		
56	C5	1	Total	Mg	0	0
			1	1		
56	13	152	Total	Mg	0	0
			152	152		
56	1J	10	Total	Mg	0	0
			10	10		
56	35	1	Total	Mg	0	0
			1	1		
56	16	15	Total	Mg	0	0
			15	15		
56	25	2	Total	Mg	0	0
			2	2		
56	M5	1	Total	Mg	0	0
			1	1		
56	21	2	Total	Mg	0	0
			2	2		
56	31	1	Total	Mg	0	0
			1	1		
56	L8	1	Total	Mg	0	0
			1	1		
56	15	1	Total	Mg	0	0
			1	1		
56	3I	2	Total	Mg	0	0
			2	2		
56	I8	2	Total	Mg	0	0
			2	2		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
56	52	1	Total 1	Mg 1	0	0
56	5E	1	Total 1	Mg 1	0	0
56	29	2	Total 2	Mg 2	0	0
56	41	2	Total 2	Mg 2	0	0
56	2K	7	Total 7	Mg 7	0	0
56	J8	1	Total 1	Mg 1	0	0
56	39	3	Total 3	Mg 3	0	0
56	1G	88	Total 88	Mg 88	0	0
56	11	2	Total 2	Mg 2	0	0
56	1H	506	Total 506	Mg 506	0	0
56	E5	1	Total 1	Mg 1	0	0
56	88	2	Total 2	Mg 2	0	0
56	49	1	Total 1	Mg 1	0	0
56	14	426	Total 426	Mg 426	0	0
56	78	2	Total 2	Mg 2	0	0
56	6A	1	Total 1	Mg 1	0	0
56	4K	1	Total 1	Mg 1	0	0
56	G8	1	Total 1	Mg 1	0	0
56	2L	3	Total 3	Mg 3	0	0

- Molecule 57 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
57	32	1	Total 1	Zn 1	0	0
57	3E	1	Total 1	Zn 1	0	0
57	5I	1	Total 1	Zn 1	0	0
57	5A	1	Total 1	Zn 1	0	0
57	G8	1	Total 1	Zn 1	0	0
57	C5	1	Total 1	Zn 1	0	0

- Molecule 58 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
58	13	164	Total 164	O 164	0	0
58	3E	1	Total 1	O 1	0	0
58	4E	1	Total 1	O 1	0	0
58	1I	1	Total 1	O 1	0	0
58	3I	1	Total 1	O 1	0	0
58	5I	1	Total 1	O 1	0	0
58	6I	1	Total 1	O 1	0	0
58	7I	1	Total 1	O 1	0	0
58	4K	1	Total 1	O 1	0	0
58	1H	920	Total 920	O 920	0	0
58	16	6	Total 6	O 6	0	0
58	11	9	Total 9	O 9	0	0
58	21	5	Total 5	O 5	0	0
58	31	7	Total 7	O 7	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
58	78	6	Total 6	O 6	0	0
58	B8	1	Total 1	O 1	0	0
58	C8	2	Total 2	O 2	0	0
58	E8	1	Total 1	O 1	0	0
58	F8	1	Total 1	O 1	0	0
58	G8	3	Total 3	O 3	0	0
58	J8	1	Total 1	O 1	0	0
58	L8	3	Total 3	O 3	0	0
58	P8	1	Total 1	O 1	0	0
58	Q8	2	Total 2	O 2	0	0
58	1G	64	Total 64	O 64	0	0
58	32	1	Total 1	O 1	0	0
58	7A	1	Total 1	O 1	0	0
58	14	543	Total 543	O 543	0	0
58	1J	18	Total 18	O 18	0	0
58	19	3	Total 3	O 3	0	0
58	29	3	Total 3	O 3	0	0
58	39	8	Total 8	O 8	0	0
58	15	1	Total 1	O 1	0	0
58	35	1	Total 1	O 1	0	0
58	85	1	Total 1	O 1	0	0

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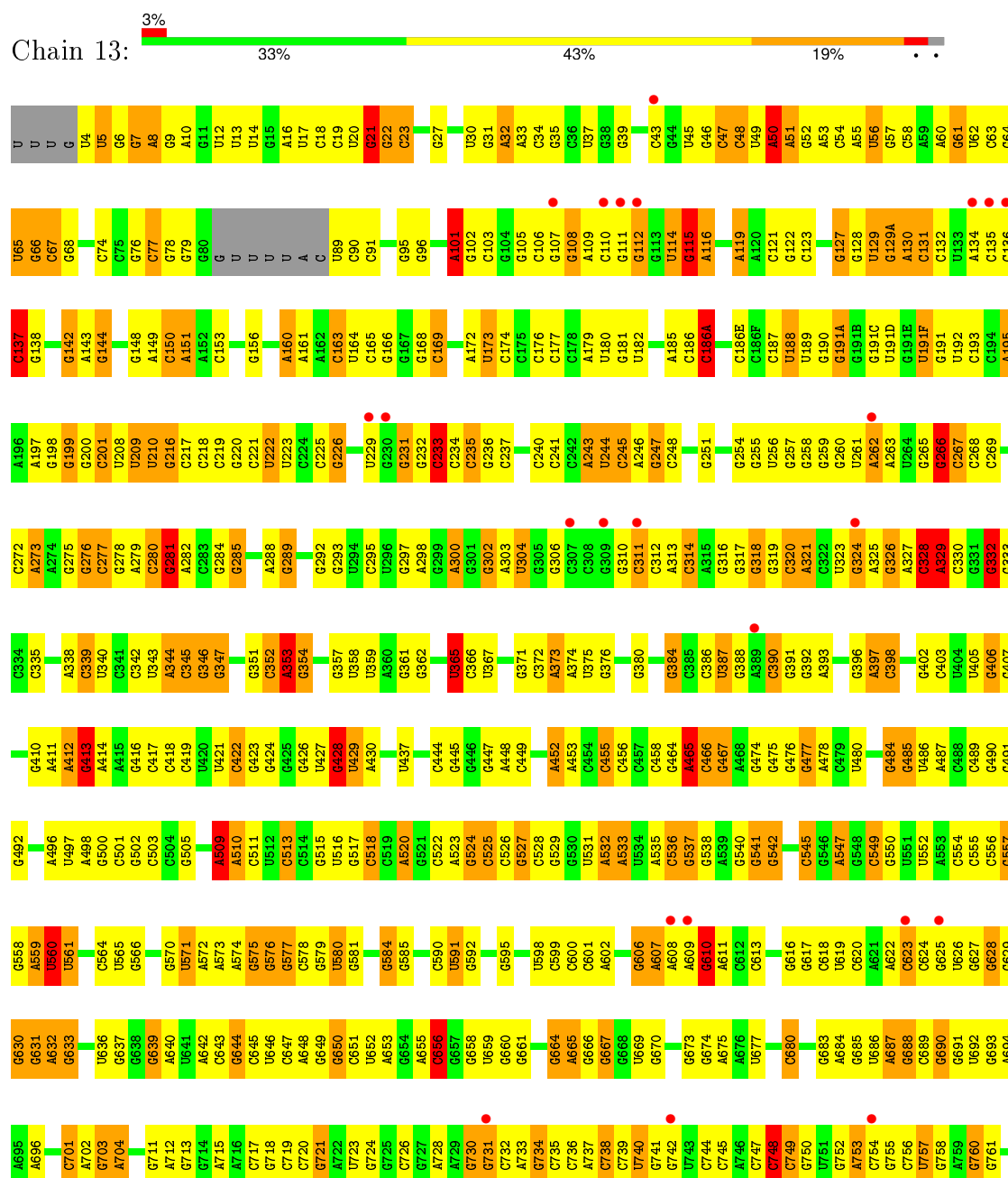
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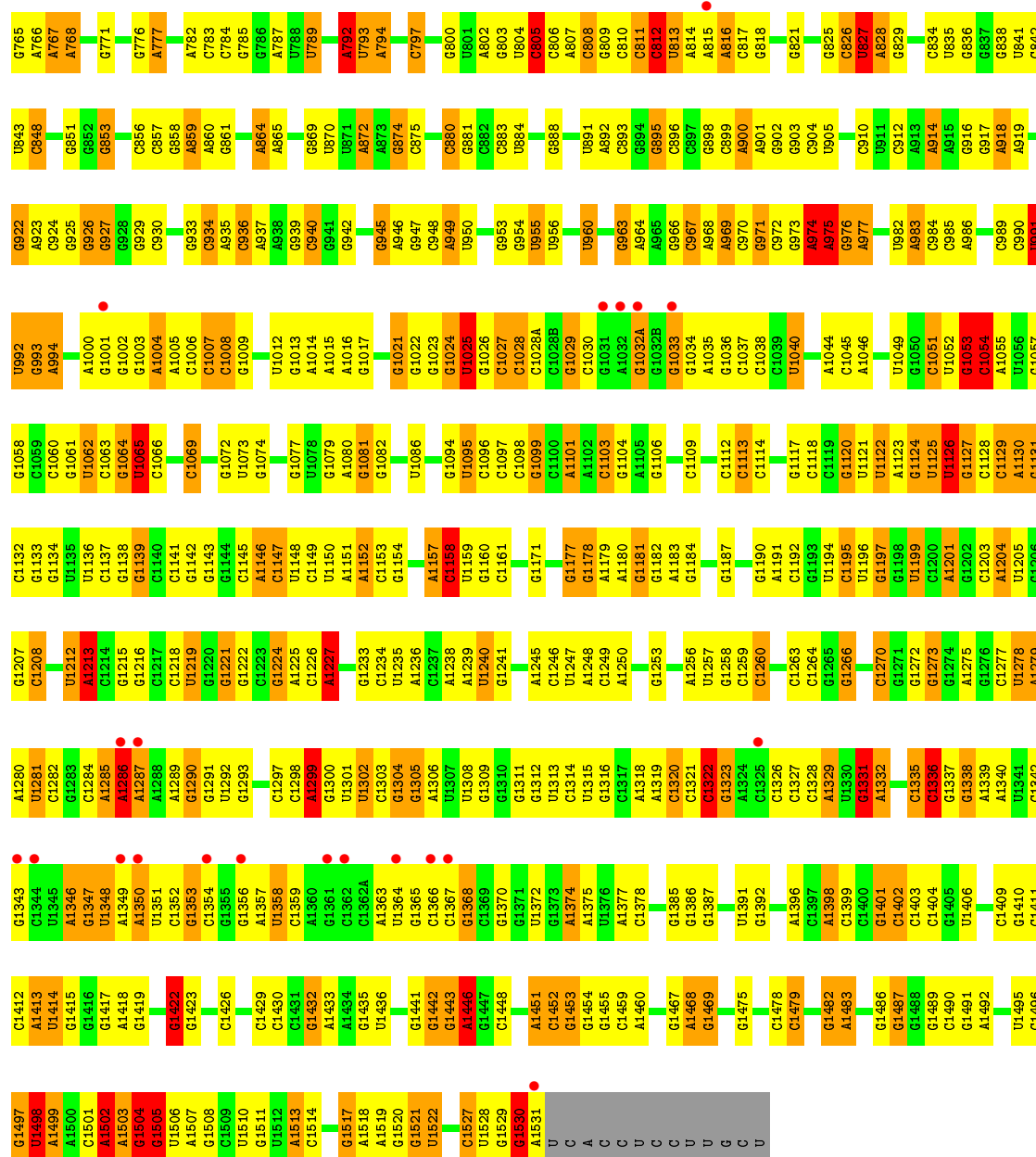
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
58	L5	2	Total	O	0	0
			2	2		

3 Residue-property plots

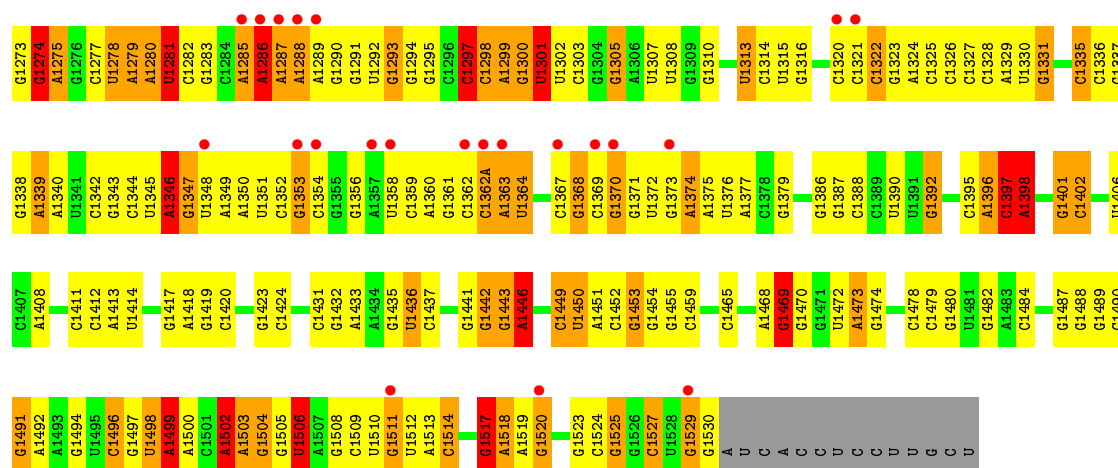
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($\text{RSRZ} > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: 16S ribosomal RNA

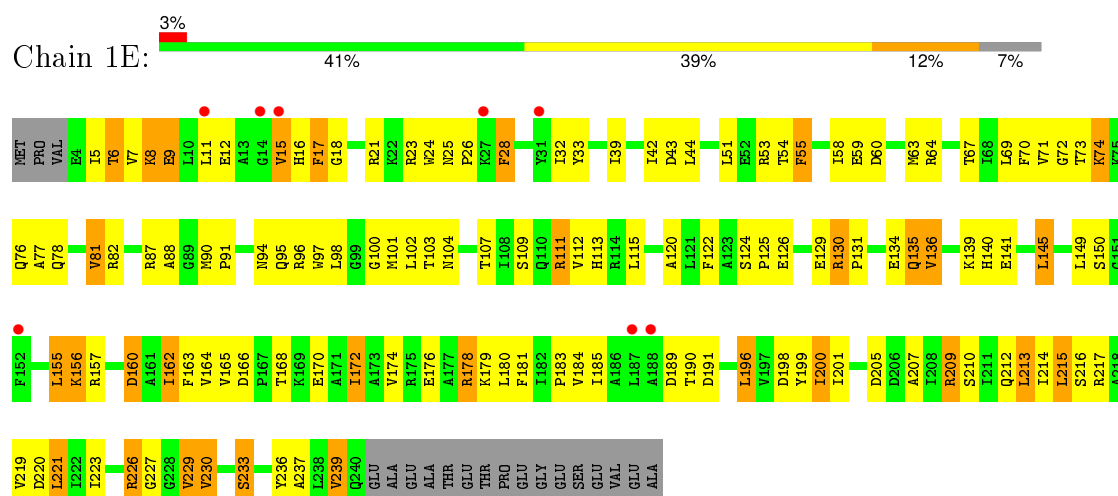




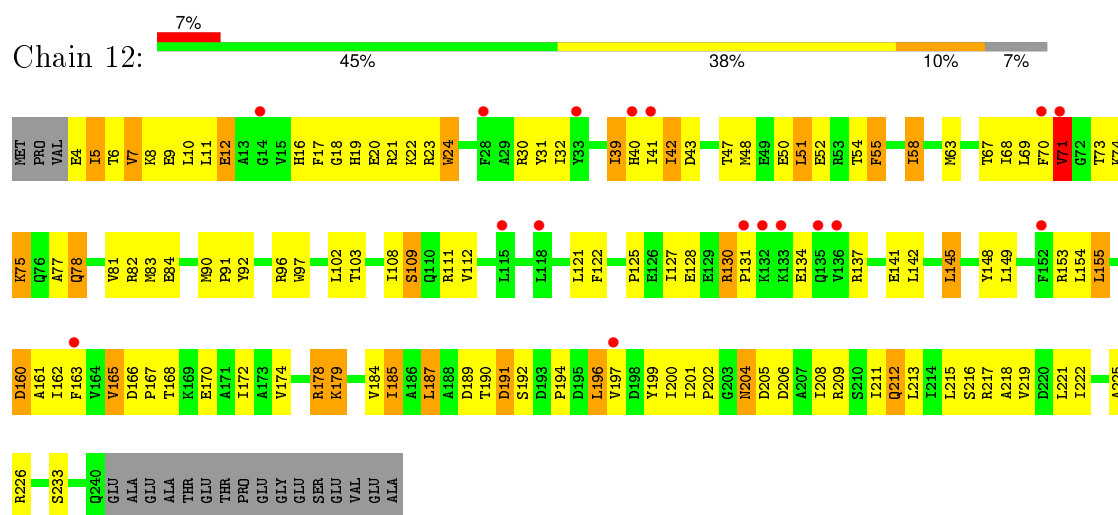
G1207	G1143	G1074	A1015	U950	G880	G811	G731	G660	G594	G521	C442	U367	G297	G216
C1208	G1144	G1078	A1016	G951	G881	C811	C735	G661	G595	C522	C443	G371	A298	C217
C1209	C1145	U1079	G1017	U952	G882	U813	C736	G662	C586	A523		G372	G299	C218
C1210	A1146	G1079	C1018	G953	C883	U813	C736	A663	G587	G524	G446	C371	A300	
U1211	A1147	A1080	C1019	G954	U884	A814	A814	G664	C588	C525		A373	G302	U222
U1212	U1148	G1081	U1020	U955	G885	A815	C738	A665	C589	C526	A448	A374	G303	U223
A1213	U1149	G1082		U956	G886	A816	C739	G666	C590	G527	C449	U375	A303	
C1214	U1150	U1083	G1024	U957	G887	C817	C739	G667	C591	C528	G450	G376	U304	G230
C1215	A1151	U1084	U1025	A958	G888	G818	G742	G668		G529	A451	G377		
	A1152	U1085	G1026	A959	G889	A819	U743		C596	G530	A452		G308	G236
	C1153	U1086	C1027	U960	G890	U820	U744		G531	U531	A453	G380	C309	C237
	G1154		U1028	U961	C893	G821	C745		G600	A532	C454			G238
U1219	G1155	G1094	C1028A	C962	G894	G822	A746		G601	A533	C455	A382	C312	
G1221	G1156	U1095	C1028B	G963		G823	A747		A602	U534	C456	A383	A313	C241
G1222	A1157	C1096	G1029	A964	G898	C824	C748		U603	A535	C457		C314	C242
C1223	G1158	C1097	C1030	A965	G899	G825	C749		G604	C536	C458	C386	A315	A243
G1224	U1159	G1098	G1031	G966	C899	C826	G750		U605	G537	C464	U387	G316	U244
A1225	G1160	C1099	A1032	C967	A900	U827	G751		G606	G538	A465	G388	G317	C245
C1226	C1161	C1100	G1032A	A968	A901	A828	C754		A607	A539	C466	A389	A246	
A1227	C1162	A1101	G1032B	A969	G904	G829	G755		A608	G540	C467	C390	G247	
C1228	C1163	A1102	G1033	C970	G830	U831				G541	A468	G391	A321	A250
		C1103	G1034	G971	U905	G760	G761		A611	G542	G474		C322	G251
U1232	G1166	G1104	A1035	C972	G906	G836	C762		C612		G475	A397	U323	
C1234	A1167	A1105	G1036	G973	A909	G837	G763		C613	C545	G476	C398	G326	G254
U1235	A1170	G1106	C1037	A974	C910	G838	C764		C614	G546	G477	C399	A327	G255
A1236	G1171	C1107	C1038	A975	U911	U841			C615	A547	A478	C400	G327	
C1237	C1172	C1109	U1040	G976	U912	C842	G769		G616	G550	C479	C403	C328	U256
A1238	G1173	A1110		A978	C912	U843	U772		G617	G551	U480	C403	C330	G257
A1239	G1174	A1111	C1043	C979	A914	U844	U773		U618	U552	C481	G406	G331	U261
U1240		C1112	A1044	U980	A915	C849	G774		U619	A553		G407	G332	A262
G1241	G1177		C1045	U981	G916	U850	G775		C620	C554	G484	A408		
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C1244	A1179	G1117	G1047	C984	A918	G852	G776		C624	C556	G486	G409	A338	C266
A1245	A1180	G1117	C1048	C984	A919	G853	A777		G625		A487	G410	C339	G267
C1246	G1181	C1118	U1049	C985	U920	G854	G778		U626	A559	G490	A411	U340	C268
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A1248	A1183	C1120	C1051	C987	G922	G856			G628	U561	A495	A414	C342	
		U1121	U1052		A923	C857	A781		C629	C562	A496	G424	G343	C272
U1250	G1186	U1122	G1053	U991	G924	G858	A782		G630	C563	U497	G425	U344	A273
A1251	G1187	A1123	C1054	U992	C924	G859	C783		G631	A564	A498	G426	C345	A274
A1252	A1188	G1124	A1055	G993	G925	A860	C784		G632	U565	G500	U427	C346	G275
G1253	G1189	U1125	U1056	A994	G926	G861	G785		G633	G566	C501	G428		G276
C1254	G1190	U1126	G1057	C995	G927	C862			G634	G567	G502	U429	G347	C277
G1255	A1191	G1127	G1058		C932	U863	U789		G637	G568	C503	G424	A349	G278
A1256	C1192	C1128	C1059	G998	G933	A864	A790		G638	C569	G504	G425	G350	G279
U1257	G1193	C1129	C1060	C998A	C934	A865	G791		G639	G570	G505	G426	G351	C280
G1258	U1194	A1130	G1061	U999	C935	C866	A792		A640	U571	G506	U427	C352	C281
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C1262	G1198	U1135	U1065	G1003	C940		G800		U646	G575	C511	A431	A356	G285
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	C1200	A1067	A1005	G942	G942	A873	A802		A648	G577	C513	C433	U358	
G1266	A1201	G1068	C1006	U943	U943	G874	A803		G649	C578	C513	U434	U359	
C1267	G1202	C1137	C1069	G944	G944	G875	U723		G650	G579	C436	C435	A360	G289
A1268	C1203	G1138	U1070			G876	G724		C651	U580	U516	C436	G361	C290
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C1270	G1206	C1140	U1073	G1013	C948	G878	A728		A653	G582	C518	G438	A363	G293
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- Molecule 2: 30S ribosomal protein S2

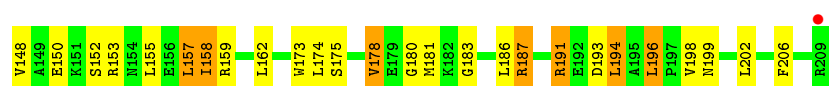


- Molecule 2: 30S ribosomal protein S2

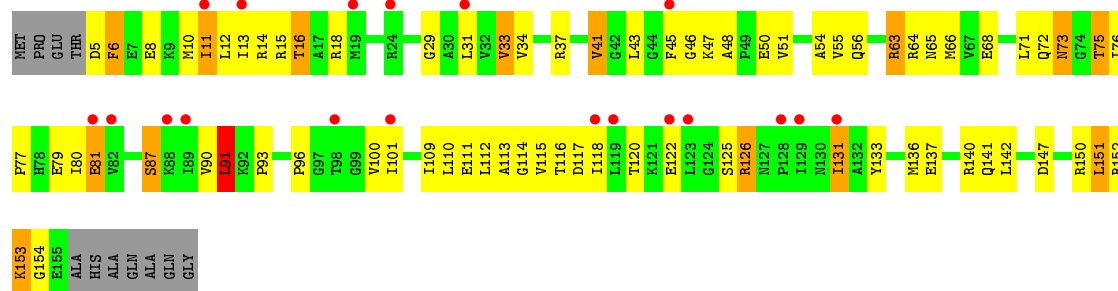


- Molecule 3: 30S ribosomal protein S3

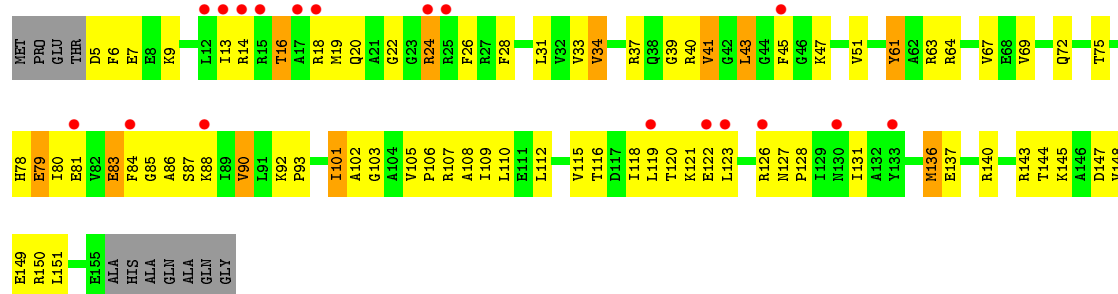




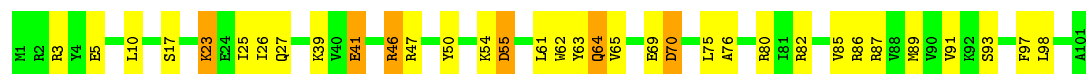
- Molecule 5: 30S ribosomal protein S5



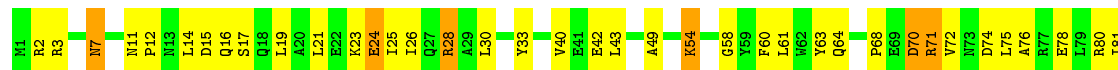
- Molecule 5: 30S ribosomal protein S5



- Molecule 6: 30S ribosomal protein S6

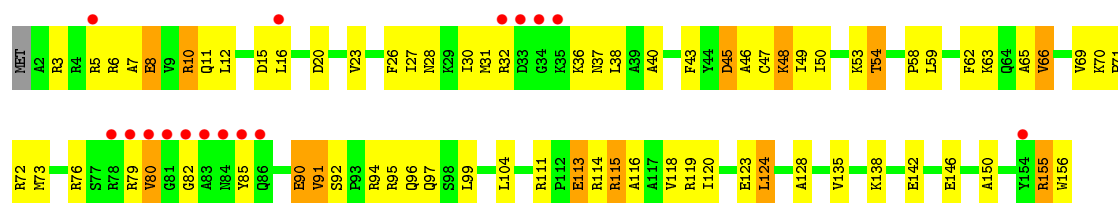


- Molecule 6: 30S ribosomal protein S6

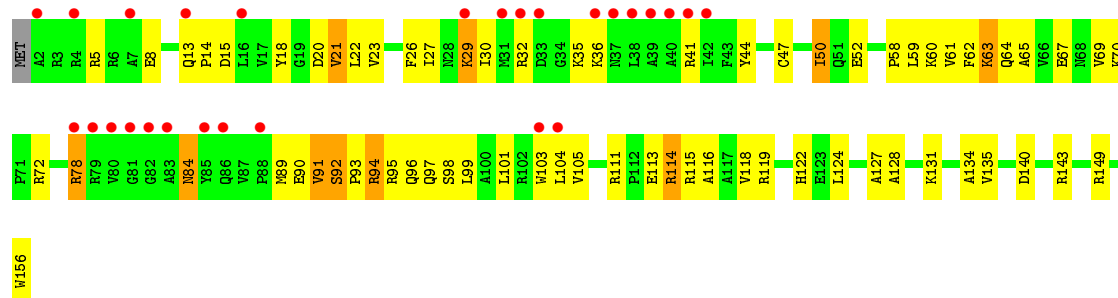


- Molecule 7: 30S ribosomal protein S7

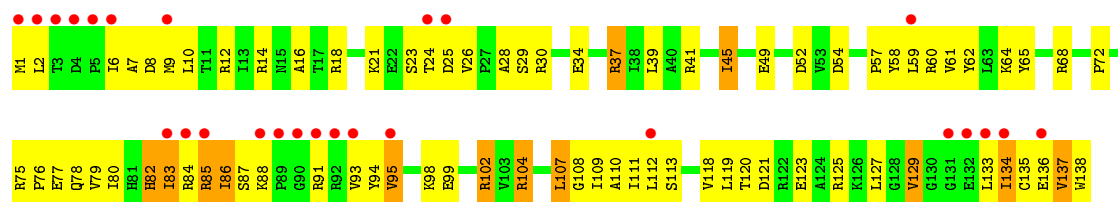




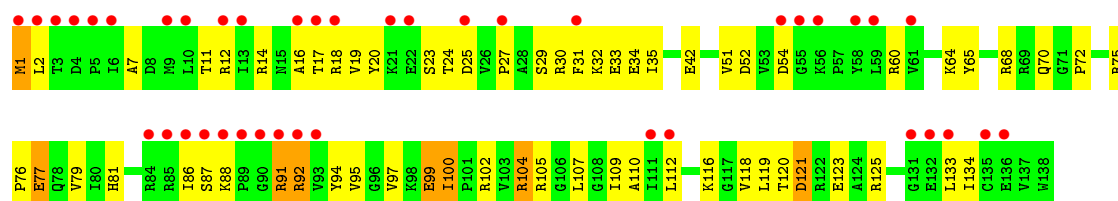
• Molecule 7: 30S ribosomal protein S7



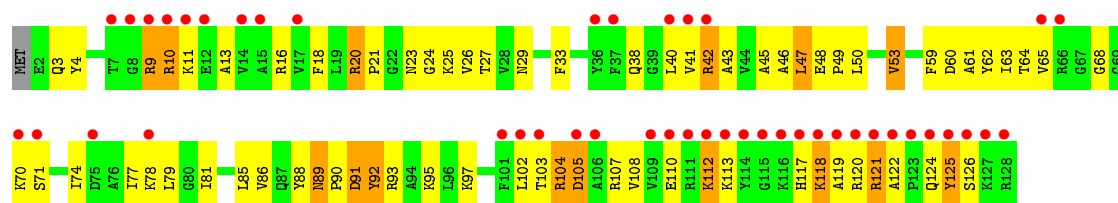
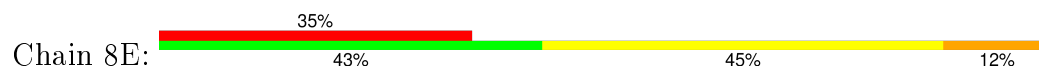
• Molecule 8: 30S ribosomal protein S8



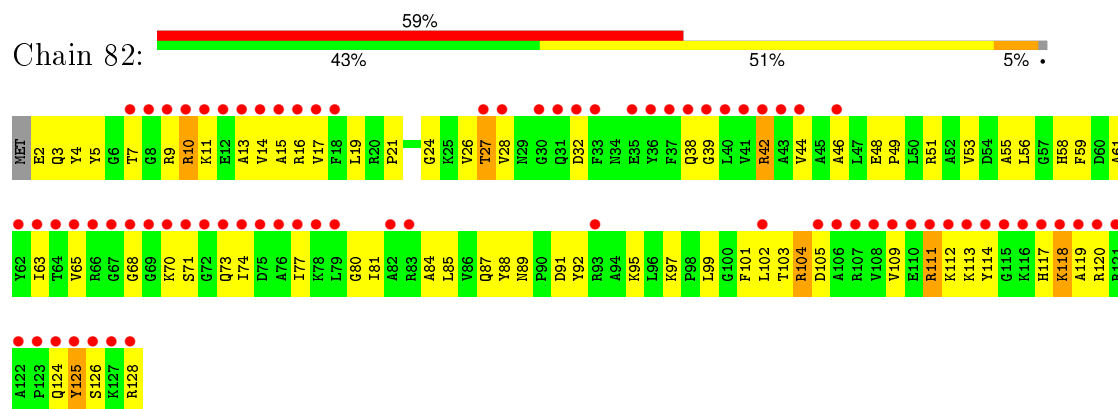
• Molecule 8: 30S ribosomal protein S8



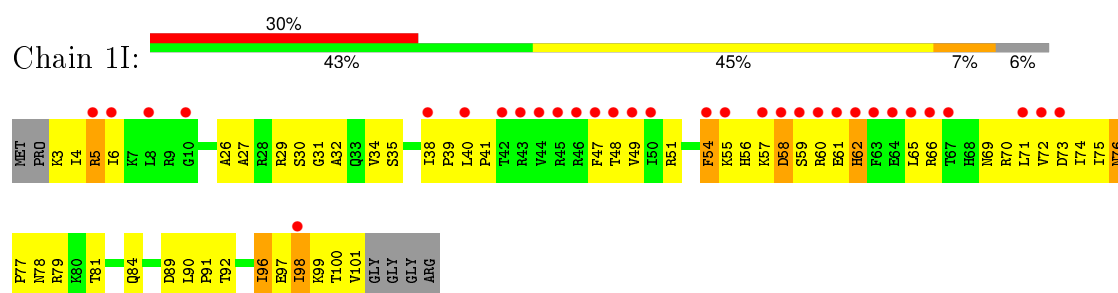
• Molecule 9: 30S ribosomal protein S9



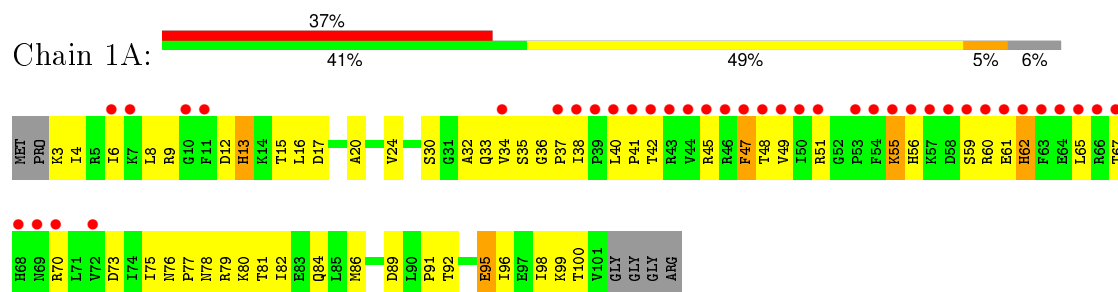
- Molecule 9: 30S ribosomal protein S9



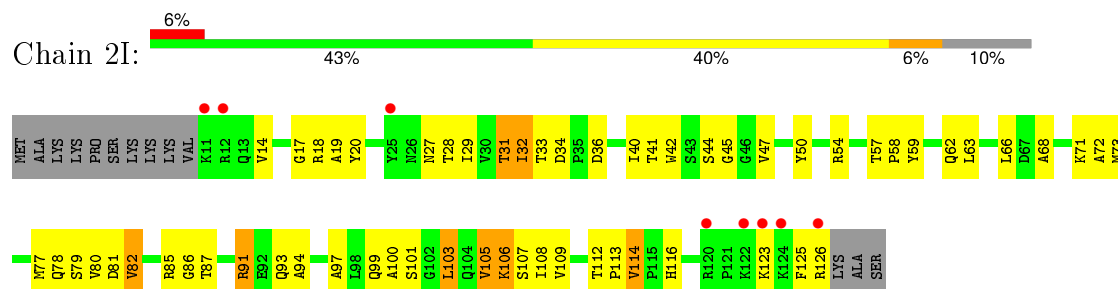
- Molecule 10: 30S ribosomal protein S10



- Molecule 10: 30S ribosomal protein S10

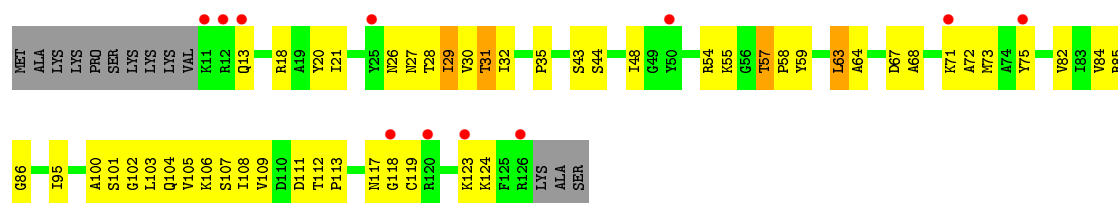


- Molecule 11: 30S ribosomal protein S11

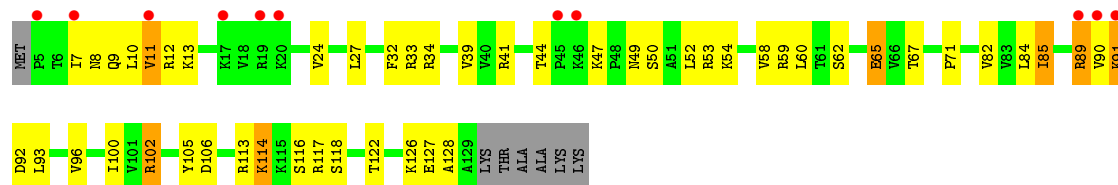


- Molecule 11: 30S ribosomal protein S11

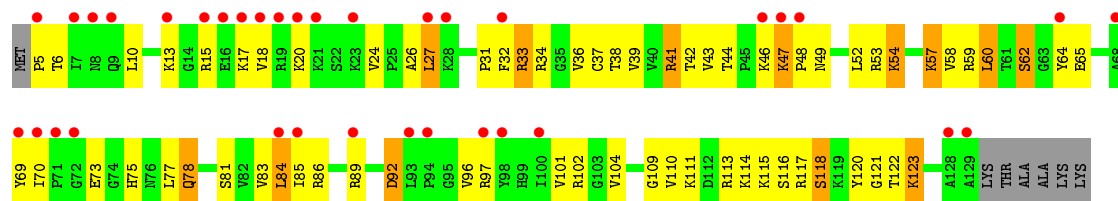
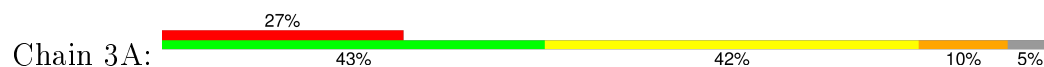




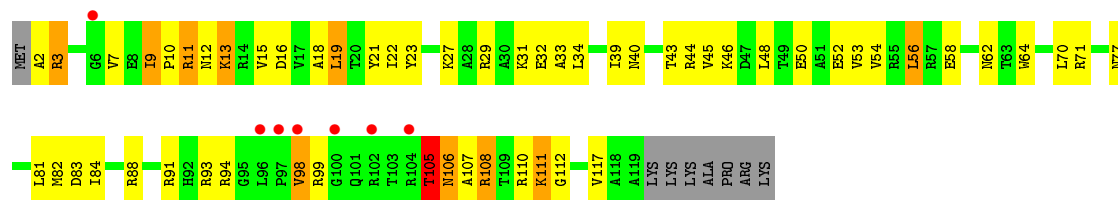
- Molecule 12: 30S ribosomal protein S12



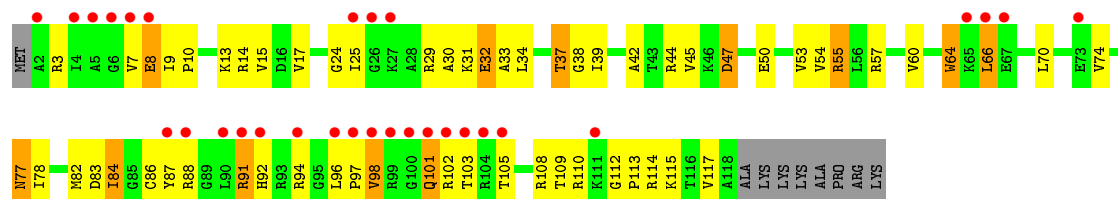
- Molecule 12: 30S ribosomal protein S12



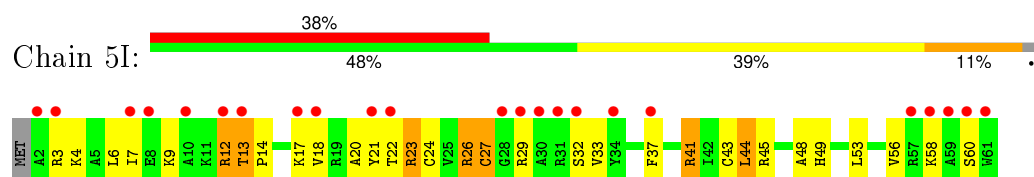
- Molecule 13: 30S ribosomal protein S13



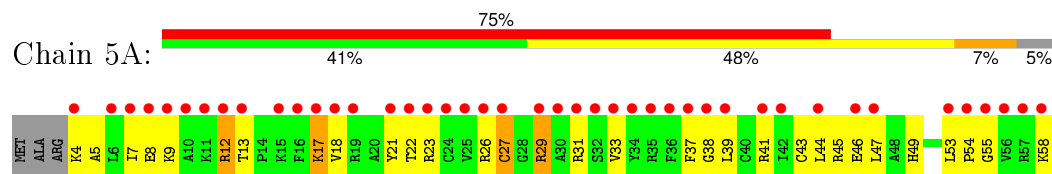
- Molecule 13: 30S ribosomal protein S13



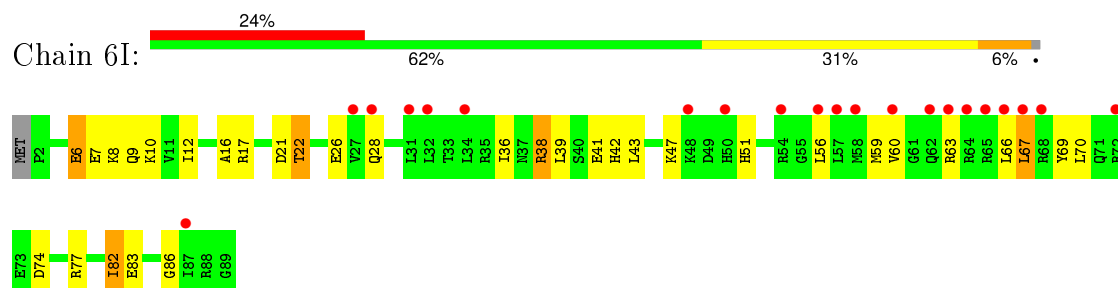
- Molecule 14: 30S ribosomal protein S14 type Z



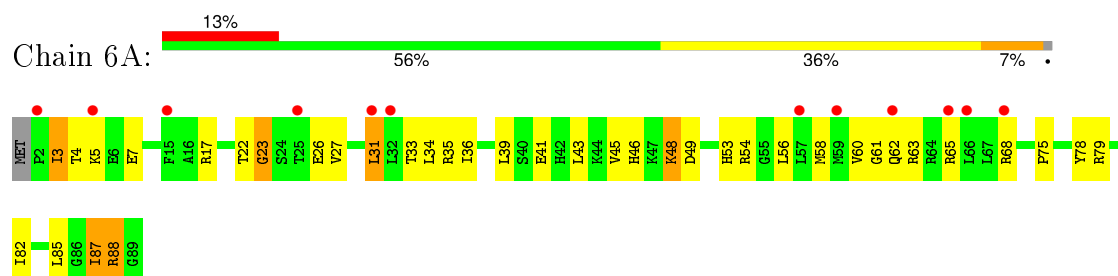
- Molecule 14: 30S ribosomal protein S14 type Z



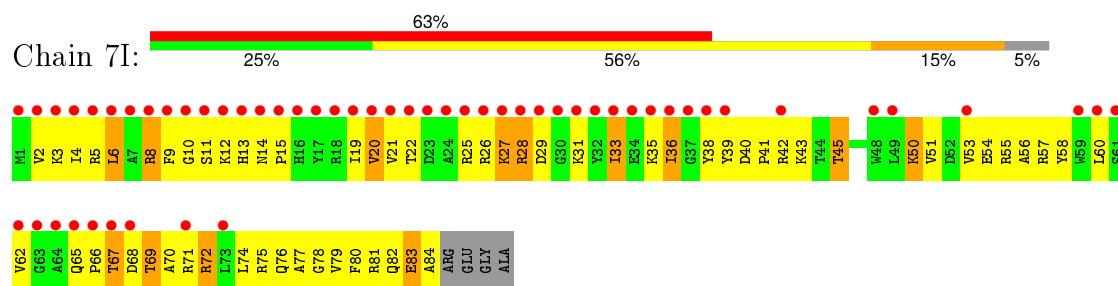
- Molecule 15: 30S ribosomal protein S15



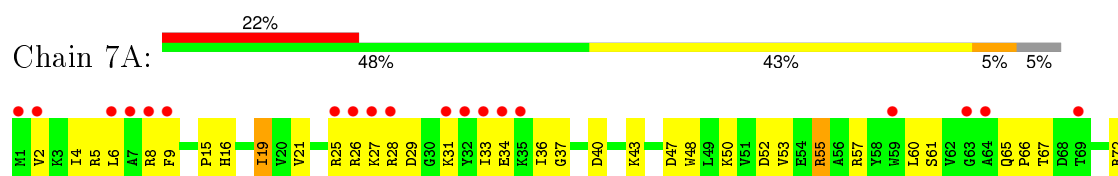
- Molecule 15: 30S ribosomal protein S15



- Molecule 16: 30S ribosomal protein S16

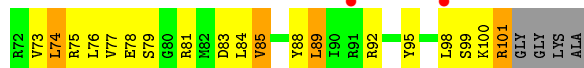
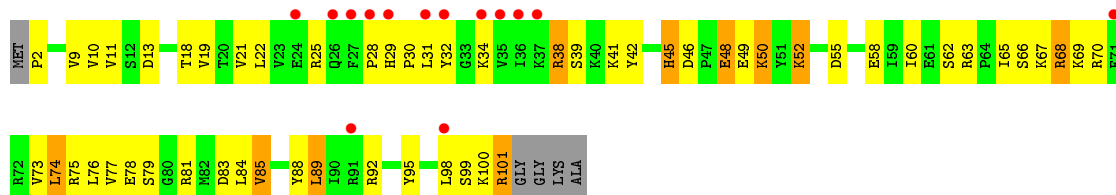


- Molecule 16: 30S ribosomal protein S16

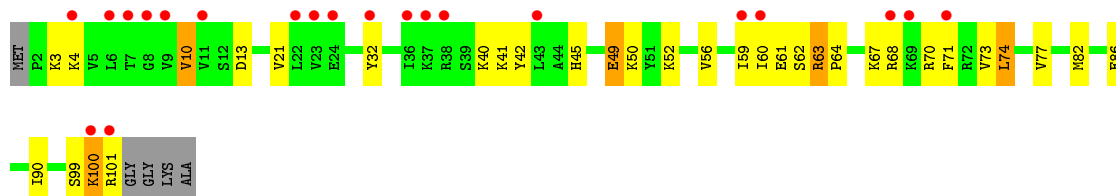




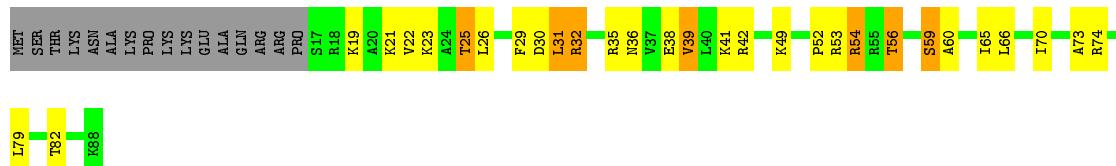
- Molecule 17: 30S ribosomal protein S17



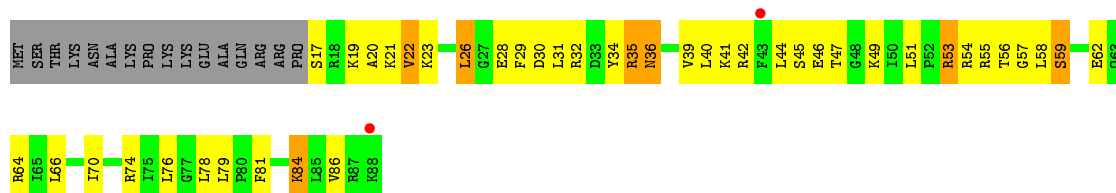
- Molecule 17: 30S ribosomal protein S17



- Molecule 18: 30S ribosomal protein S18

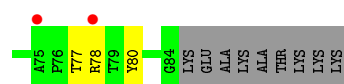


- Molecule 18: 30S ribosomal protein S18

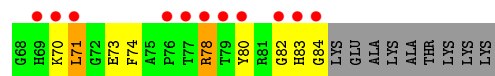
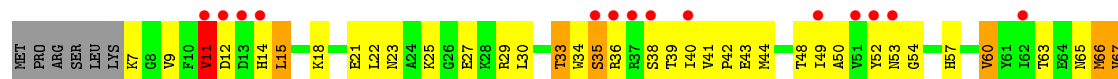


- Molecule 19: 30S ribosomal protein S19





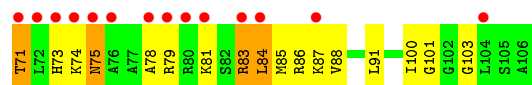
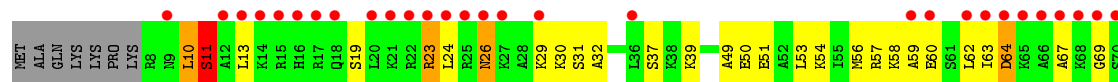
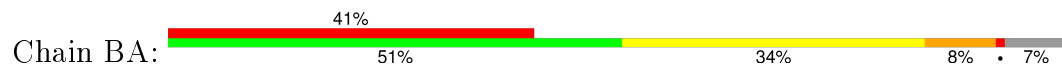
- Molecule 19: 30S ribosomal protein S19



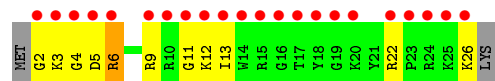
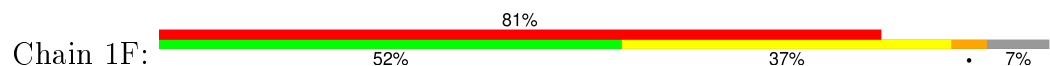
- Molecule 20: 30S ribosomal protein S20



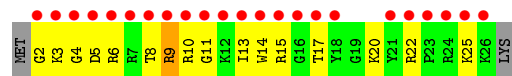
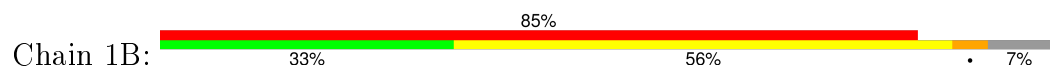
- Molecule 20: 30S ribosomal protein S20



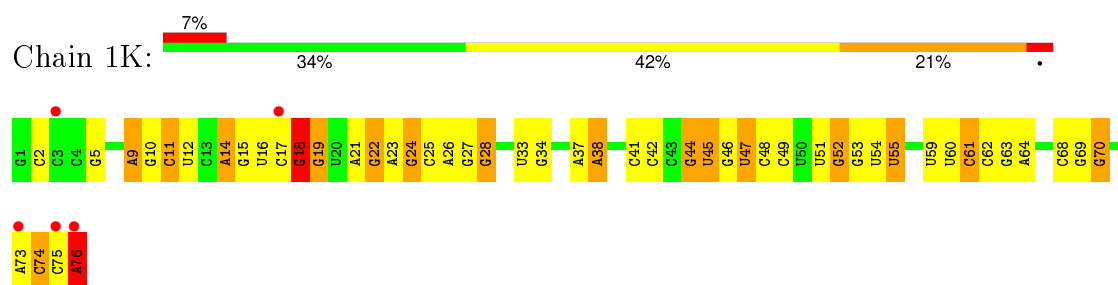
- Molecule 21: 30S ribosomal protein Thx



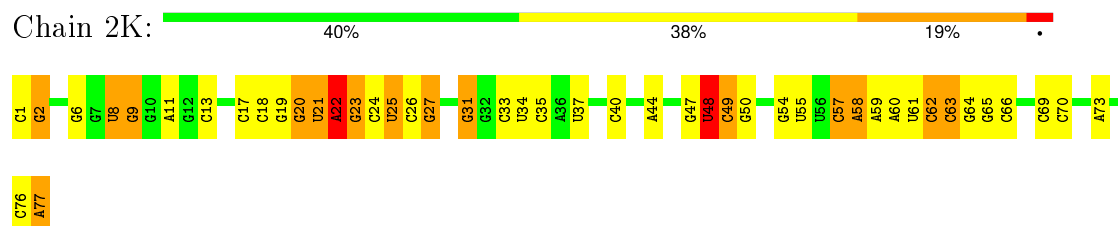
- Molecule 21: 30S ribosomal protein Thx



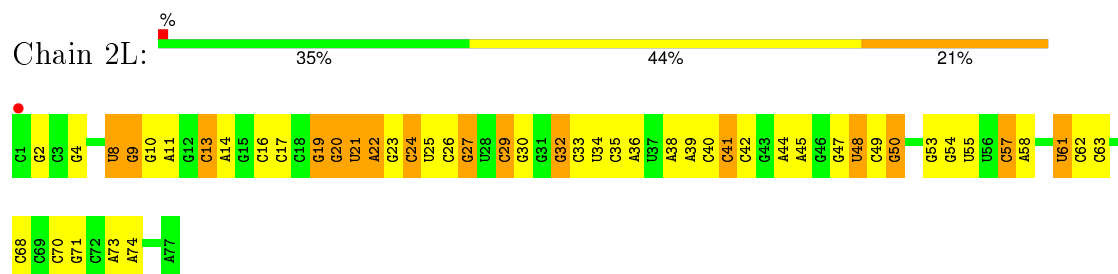
- Molecule 22: tRNA-Phe



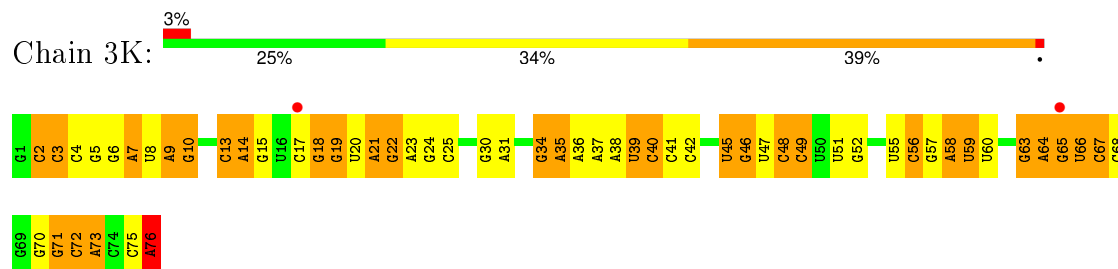
- Molecule 23: tRNA-fMet



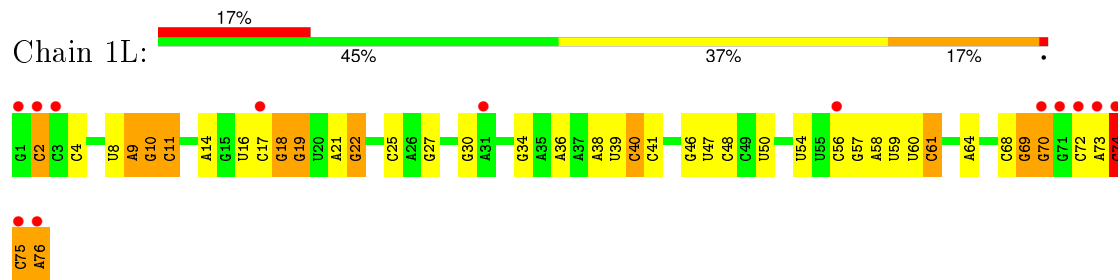
- Molecule 23: tRNA-fMet



- Molecule 24: tRNA-Phe

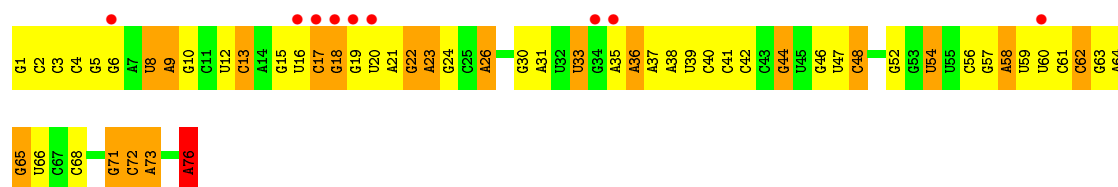


- Molecule 24: tRNA-Phe



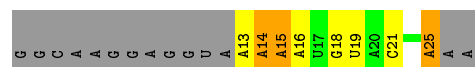
- Molecule 24: tRNA-Phe





• Molecule 25: mRNA

Chain 4K: 19% 19% 11% 52%



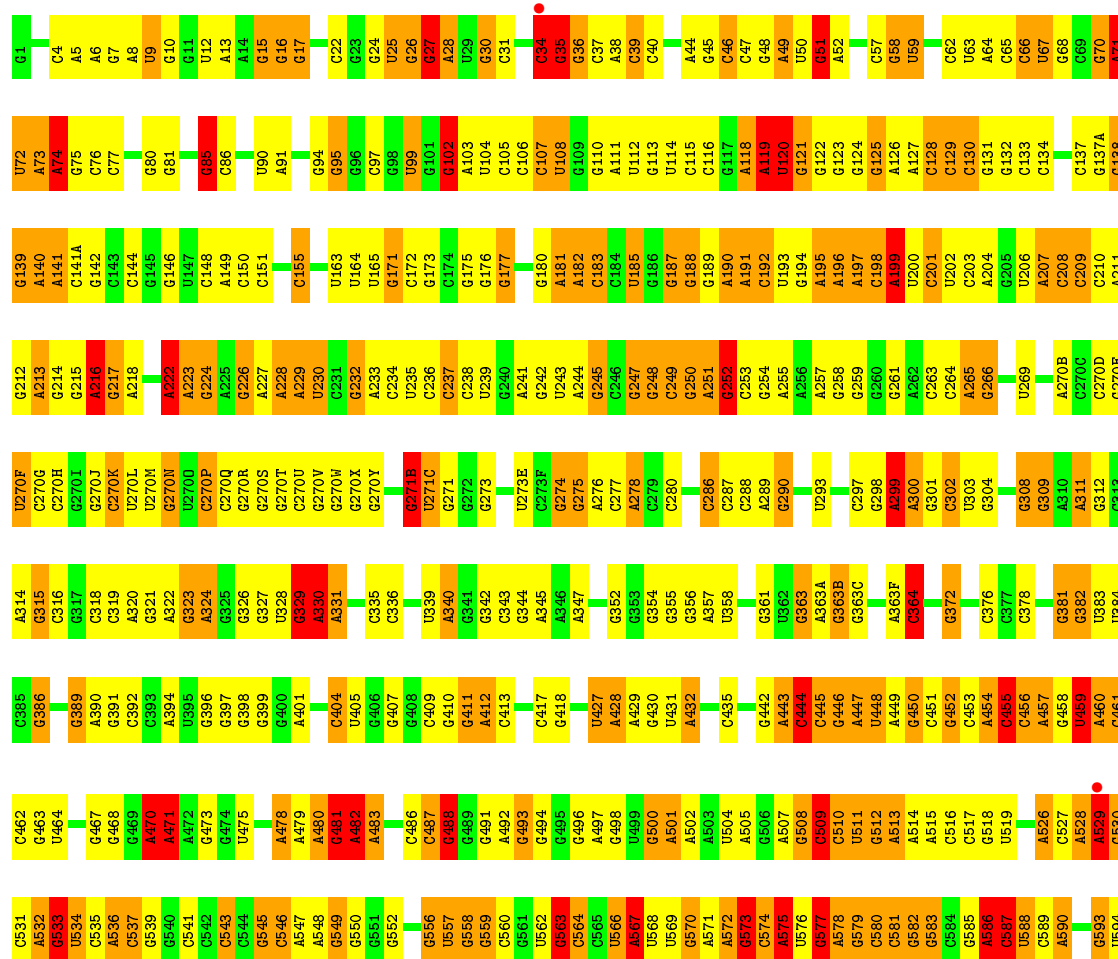
• Molecule 25: mRNA

Chain 4L: 7% 15% 11% 67%



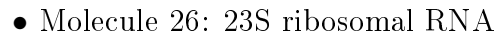
• Molecule 26: 23S ribosomal RNA

Chain 1H: 26% 42% 26% 6%

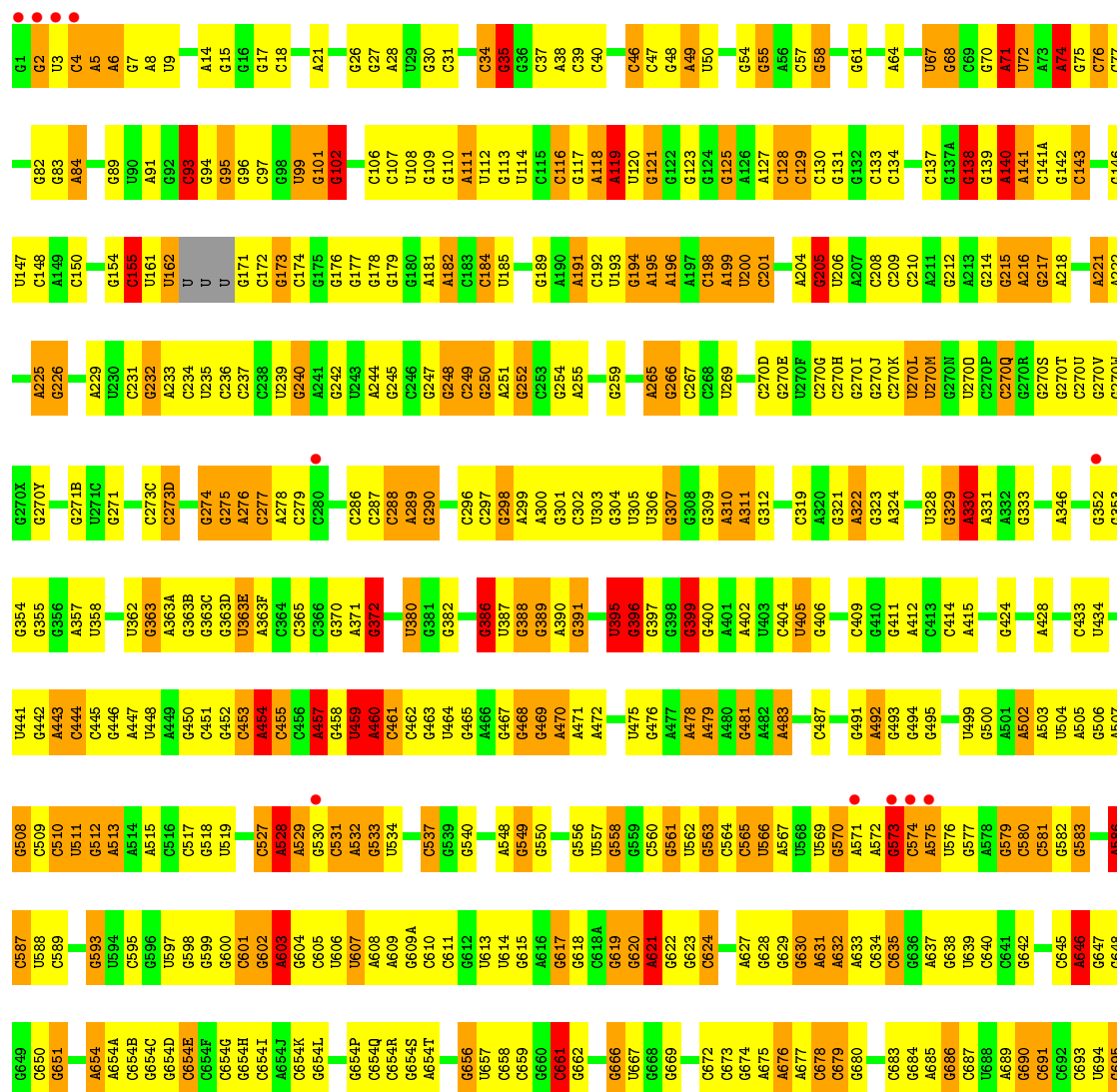




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U2604	U2530	C2453	A2327	C2393	A2267	C2193	G2055	C1986	C1925	C1836	U1775	C1690	C1624
C2605	A2531	C2456	G2328	C2394	U2268	C2194	C2056	C1989	U1926	C1837	U1776	U1692	C1625
C2606	U2457	C2457	G2329	C2395	A2269	U2197	A2057	C1990	A1927	C1838	U1777	U1693	G1626
C2607	G2458	C2458	G2330	C2396	G2270	U2198	A2058	U1993	A1928	G1839	U1778	G1694	U1629
G2608	C2543	C2459	G2331	G2397	G2271	A2199	A2059	G1997	G1929	G1842	U1779	U1695	U1630
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C2610	G2545	C2462	A2335	G2399	A2273	C2206	G2061	A2001	U1931	A1847	C1781	G1697	A1633
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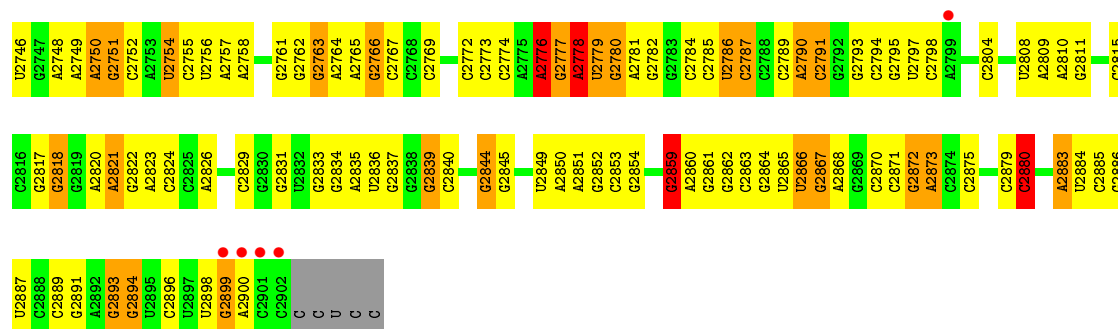


Chain 14: 2% 31% 43% 22%



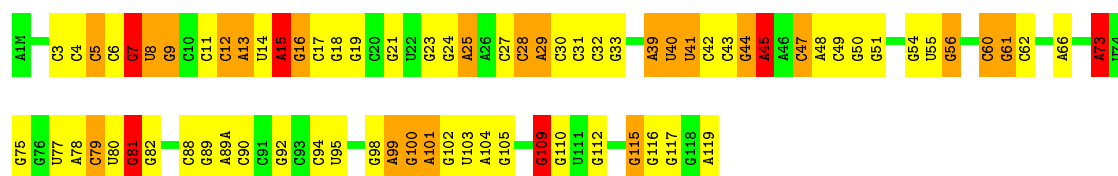
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G1642	C1505	U1576	C1506	G1436	G1371	A1307	G1244	U1167	U1097	U1032	U969	U907	G846	G776	A699
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G1653	C1513	C1586	C1513	G1445	G1380	U1316	A1253	G1176	C1104	G1039	C976	C915	G852	A784	G717
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A1701	A1494	G1568	A1494	G1429	G1364	U1300	U1300	G1235	C1158	A1086	A1020	U958	C897	G831	C766
G1702	A1495	C1569	A1495	C1430	A1365	A1301	A1301	G1236	U1159	G1089	G1022	A960	C898	U832	G767
C1636	U1431	A1570	U1431	U1431	A1366	A1302	A1302	G1237	G1160	U1090	G1024	C961	A899	G834	G768
A1637	A1432	C1432	A1432	C1432	A1367	G1303	G1303	G1238	C1161	C1092	U1026	U963	C901	A835	G770
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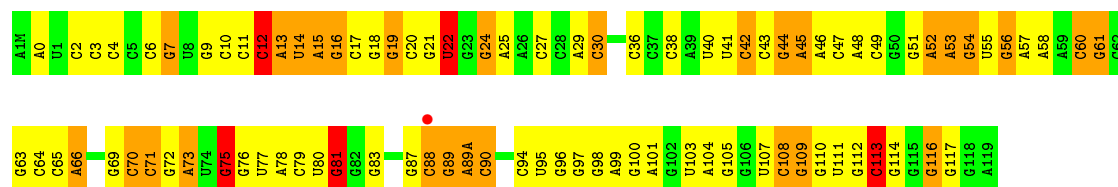
• Molecule 27: 5S ribosomal RNA

Chain 16: 38% 39% 18% 5%



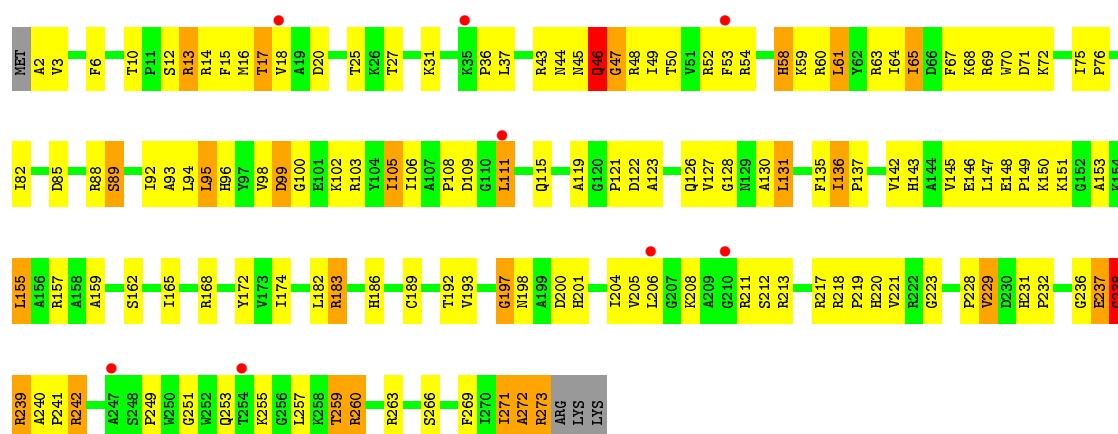
• Molecule 27: 5S ribosomal RNA

Chain 1J: 26% 47% 23% .



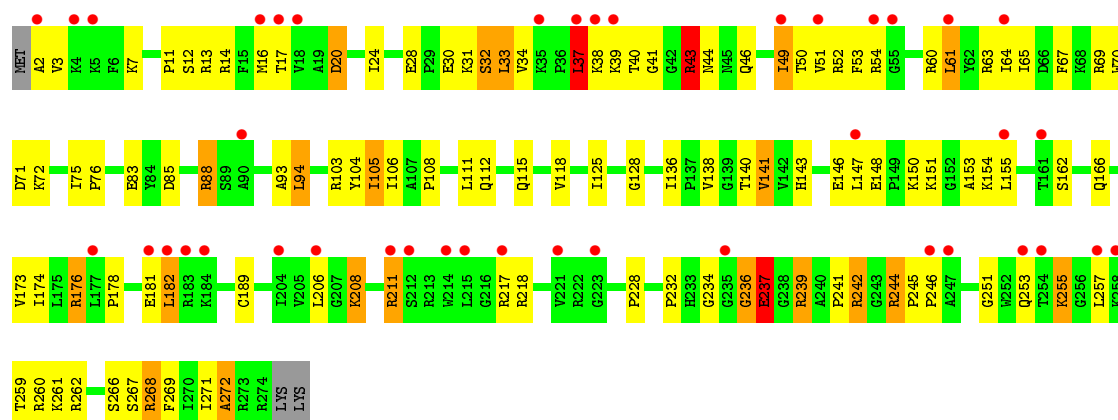
• Molecule 28: 50S ribosomal protein L2

Chain 11: 3% 48% 41% 9% ..

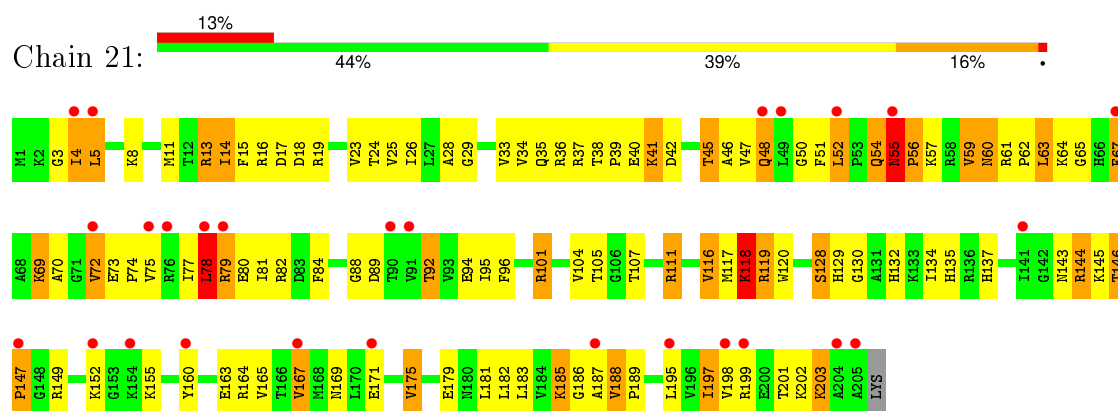


• Molecule 28: 50S ribosomal protein L2

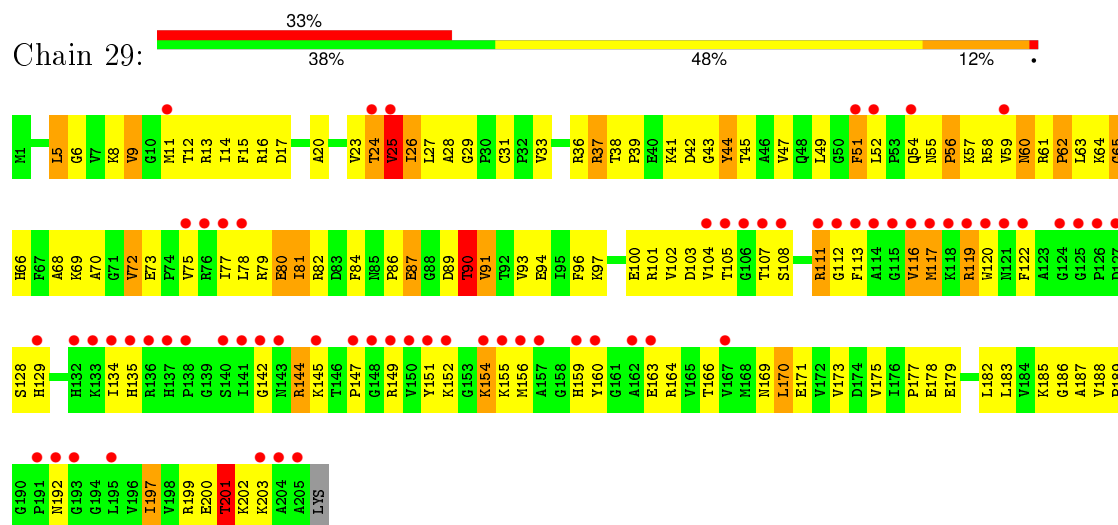
Chain 19: 15% 59% 32% 7% ..



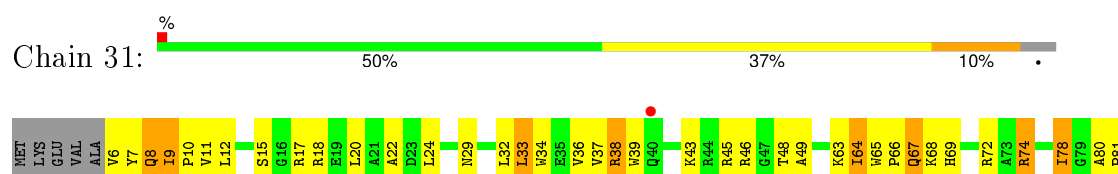
• Molecule 29: 50S ribosomal protein L3

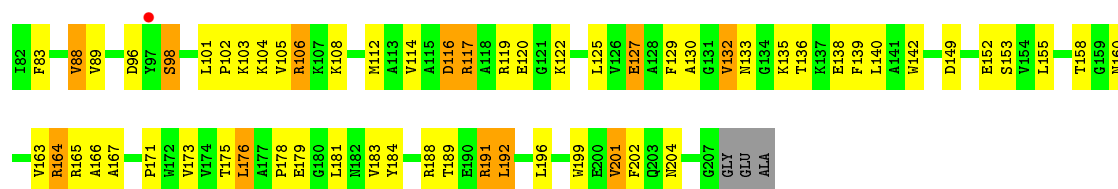


• Molecule 29: 50S ribosomal protein L3

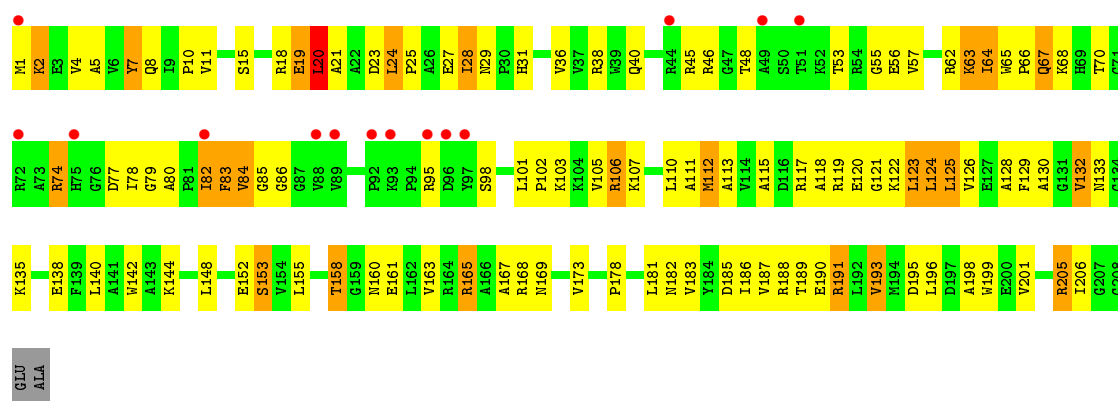


• Molecule 30: 50S ribosomal protein L4

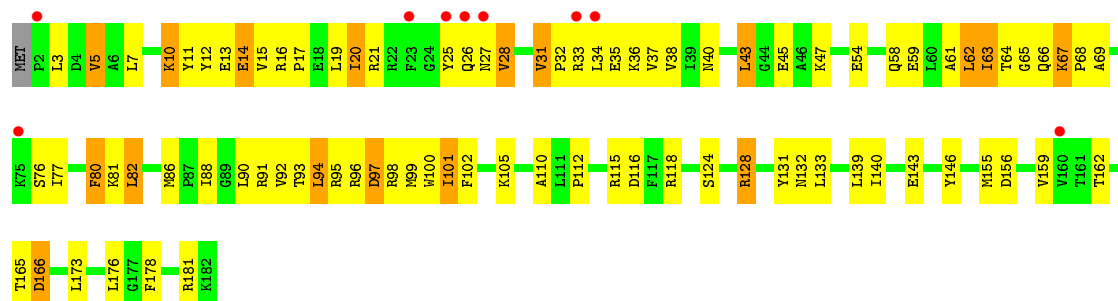




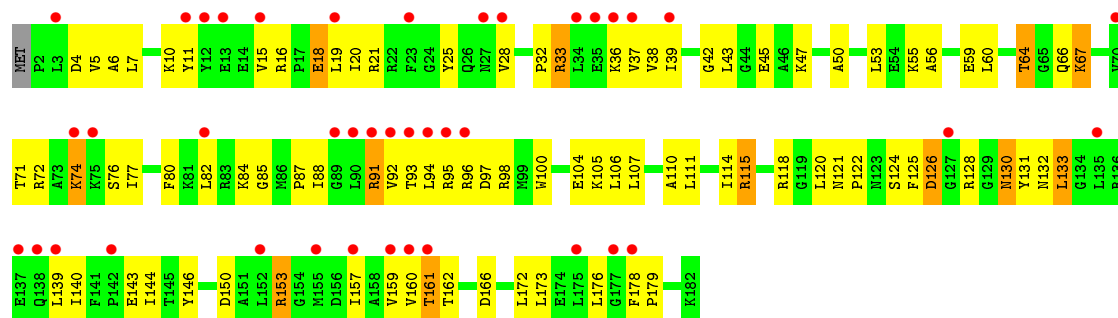
- Molecule 30: 50S ribosomal protein L4



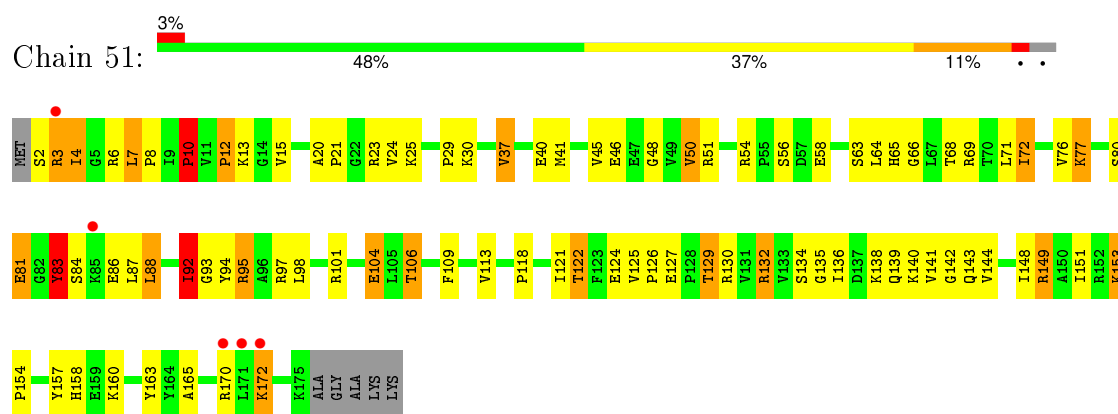
- Molecule 31: 50S ribosomal protein L5



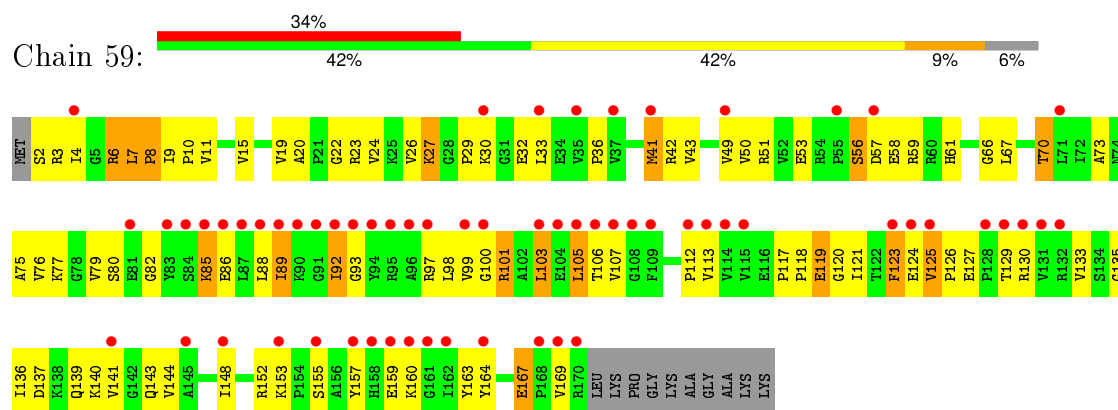
- Molecule 31: 50S ribosomal protein L5



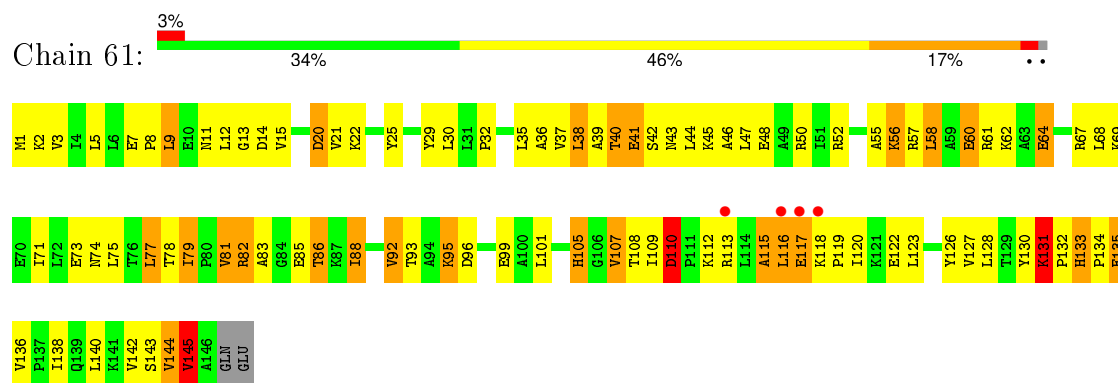
- Molecule 32: 50S ribosomal protein L6



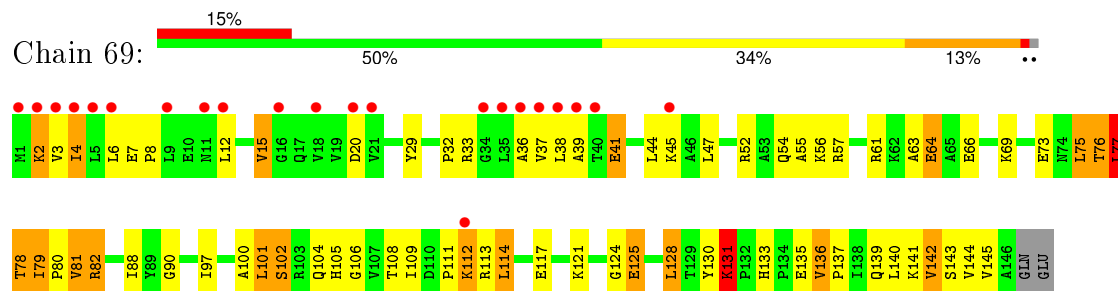
• Molecule 32: 50S ribosomal protein L6



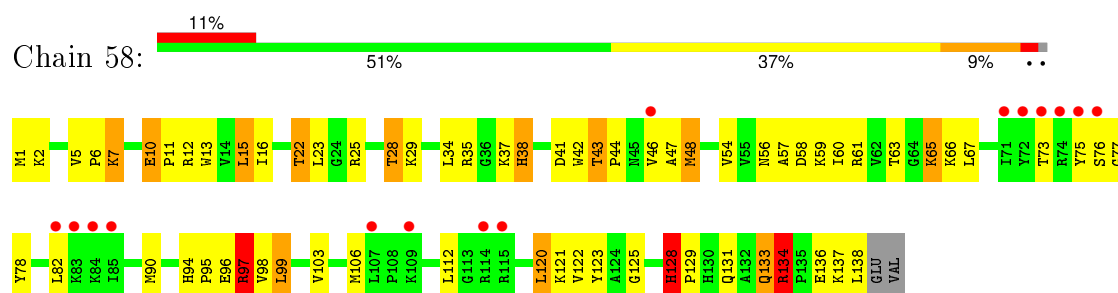
• Molecule 33: 50S ribosomal protein L9



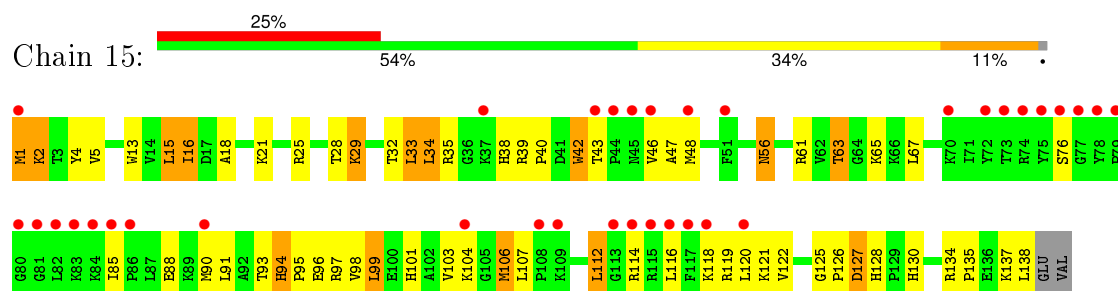
• Molecule 33: 50S ribosomal protein L9



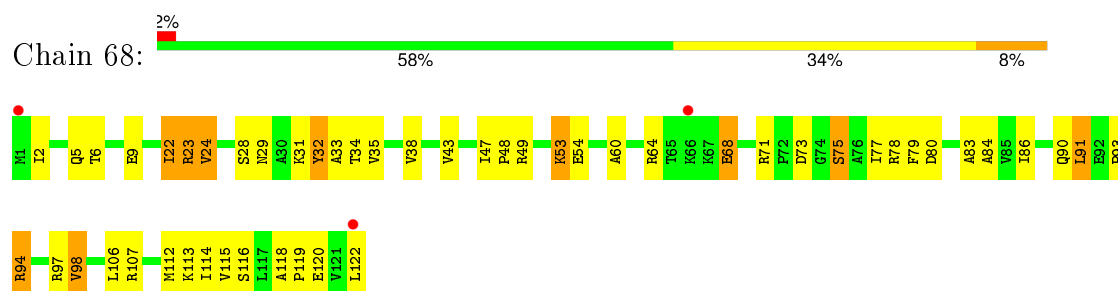
• Molecule 34: 50S ribosomal protein L13



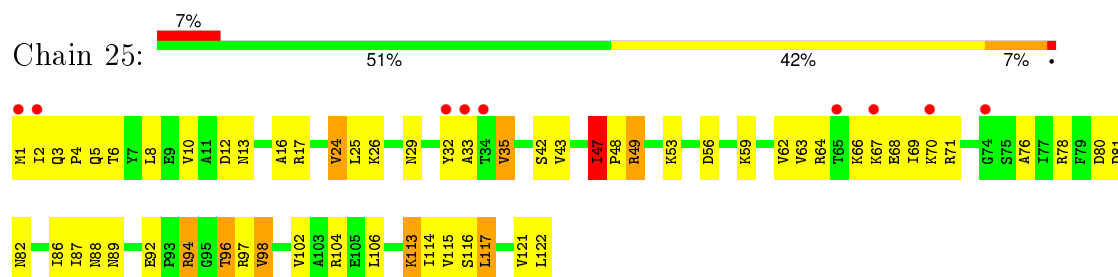
• Molecule 34: 50S ribosomal protein L13



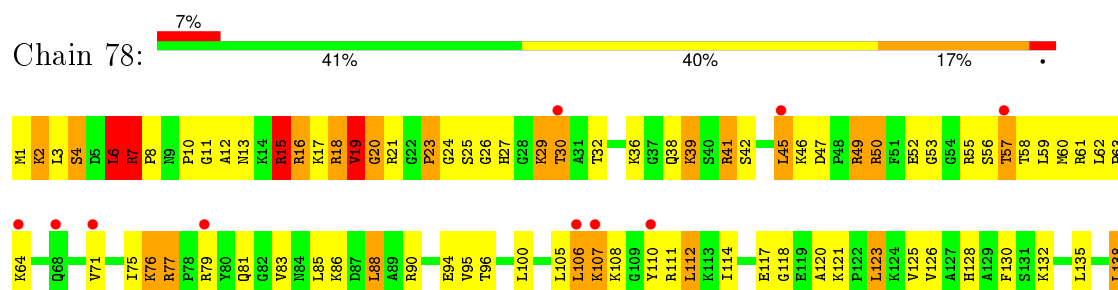
• Molecule 35: 50S ribosomal protein L14



• Molecule 35: 50S ribosomal protein L14

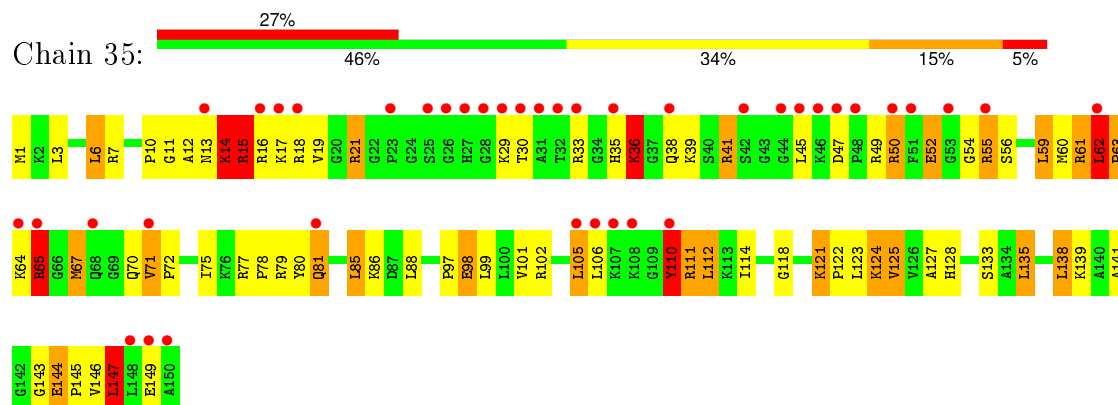


• Molecule 36: 50S ribosomal protein L15

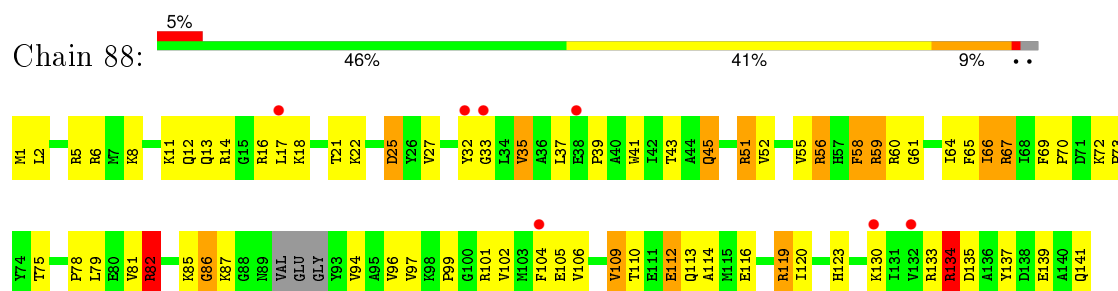




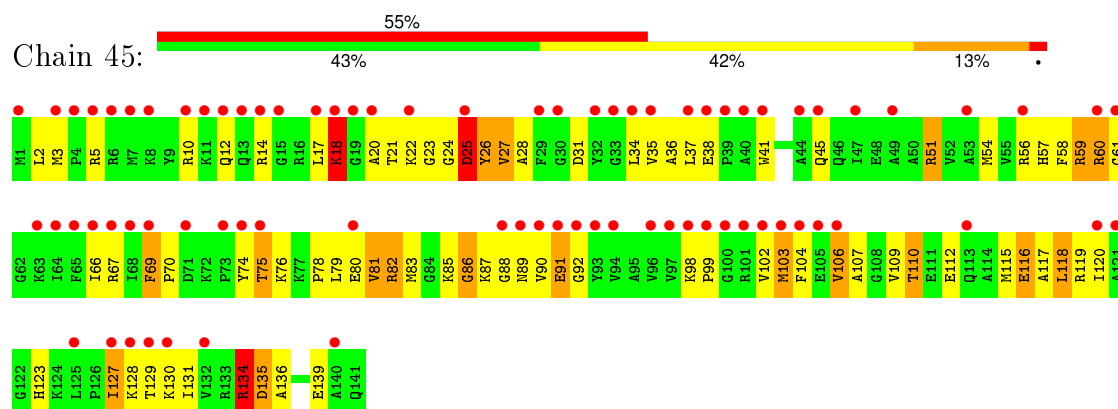
- Molecule 36: 50S ribosomal protein L15



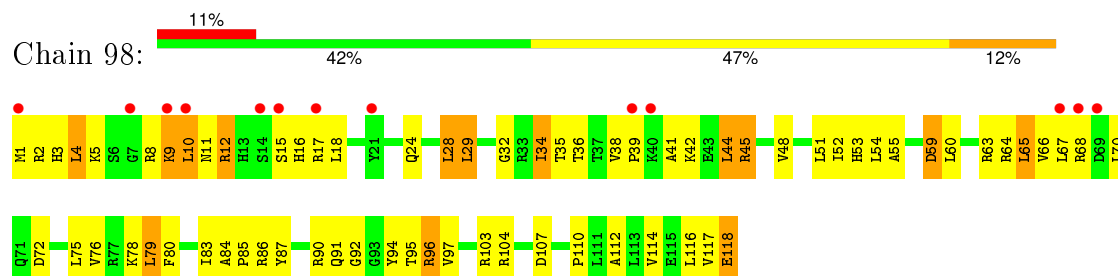
- Molecule 37: 50S ribosomal protein L16



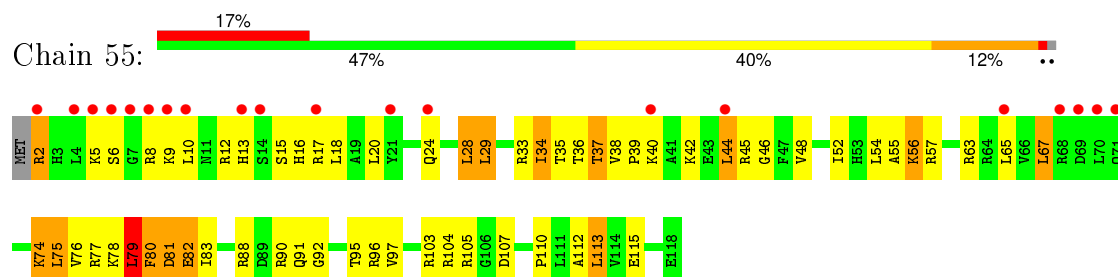
- Molecule 37: 50S ribosomal protein L16



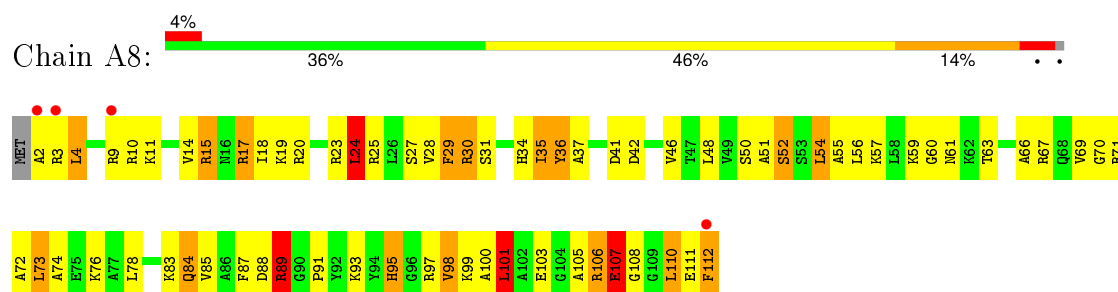
- Molecule 38: 50S ribosomal protein L17



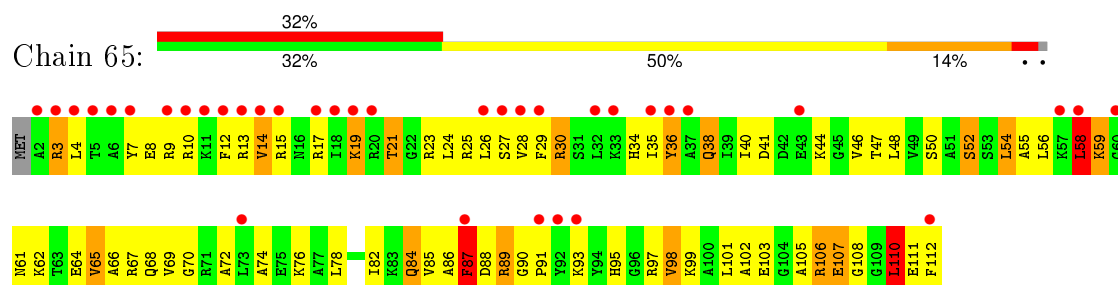
- Molecule 38: 50S ribosomal protein L17



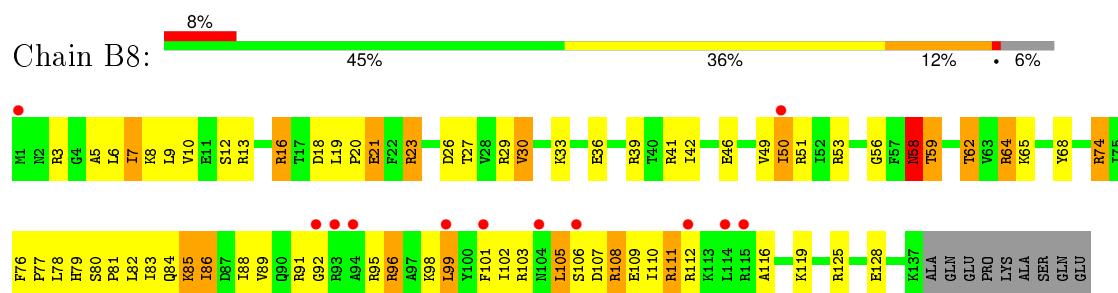
- Molecule 39: 50S ribosomal protein L18



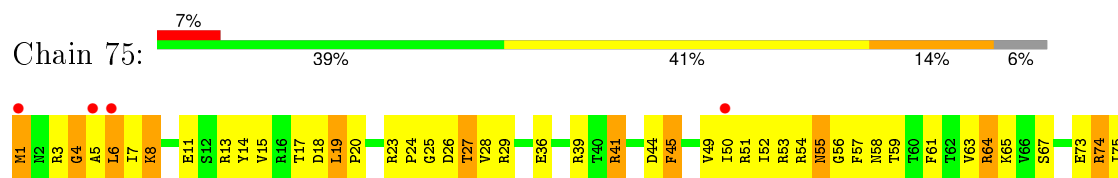
- Molecule 39: 50S ribosomal protein L18



- Molecule 40: 50S ribosomal protein L19

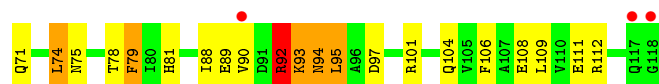


- Molecule 40: 50S ribosomal protein L19

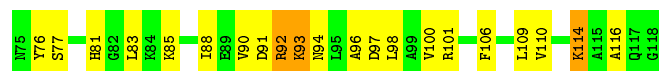




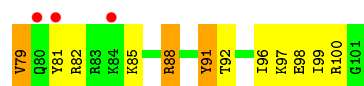
- | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
|-----|----|----|----|----|--|----|----|-----|-----|-----|-----|-----|-----|-----|-----|-----|--|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|--|-----|--|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|--|-----|-----|-----|-----|
| MET | P2 | R3 | A4 | K5 | | V8 | V9 | R10 | R11 | R12 | R13 | H14 | K15 | K16 | I17 | L18 | | G22 | G23 | Y24 | Y25 | G26 | L27 | R28 | S29 | K30 | S31 | F32 | R33 | K34 | A35 | R36 | E37 | T38 | L39 | F40 | A41 | A42 | G43 | | M44 | | H49 | R50 | K51 | R52 | K53 | K54 | R55 | D56 | F57 | R58 | R59 | L60 | M61 | | A67 | A68 | G69 | F70 |
|-----|----|----|----|----|--|----|----|-----|-----|-----|-----|-----|-----|-----|-----|-----|--|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|--|-----|--|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|--|-----|-----|-----|-----|



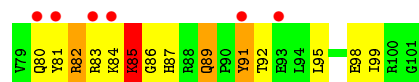
-
- | Cluster | Number of Genes |
|---------|-----------------|
| MET | 1 |
| P2 | 1 |
| P3 | 1 |
| A4 | 1 |
| K5 | 1 |
| T6 | 1 |
| G7 | 1 |
| V8 | 1 |
| R12 | 1 |
| K13 | 1 |
| H14 | 1 |
| K15 | 1 |
| K16 | 1 |
| I17 | 1 |
| L18 | 1 |
| K19 | 1 |
| L20 | 1 |
| A21 | 1 |
| K22 | 1 |
| W25 | 1 |
| R28 | 1 |
| S29 | 1 |
| K30 | 1 |
| S31 | 1 |
| F32 | 1 |
| R33 | 1 |
| K34 | 1 |
| T38 | 1 |
| L39 | 1 |
| F40 | 1 |
| G43 | 1 |
| N44 | 1 |
| Y45 | 1 |
| A46 | 1 |
| Y47 | 1 |
| A48 | 1 |
| H49 | 1 |
| R50 | 1 |
| K51 | 1 |
| R52 | 1 |
| R53 | 1 |
| K54 | 1 |
| R55 | 1 |
| D56 | 1 |
| F57 | 1 |
| R58 | 1 |
| R59 | 1 |
| R64 | 1 |
| I65 | 1 |
| M66 | 1 |
| Q71 | 1 |
| I74 | 1 |



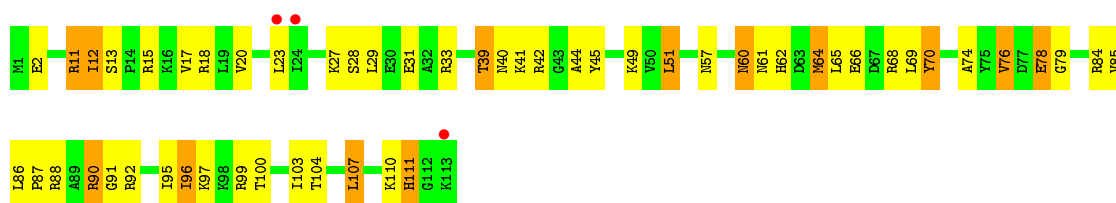
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|----|----|----|----|----|----|----|----|----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|
| M1 | F2 | A3 | I4 | V5 | K6 | I7 | G8 | G9 | K10 | Q11 | V14 | L18 | K19 | L20 | R21 | V22 | E23 | K24 | L25 | E34 | L35 | F36 | V37 | L38 | L39 | L40 | G41 | G42 | E43 | K44 | T45 | V46 | V47 | G48 | T49 | E53 | G54 | A55 | S56 | V57 | V58 | A59 | E60 | G64 | H65 | R66 | V72 | S73 | K74 | F75 | K76 | V77 | A78 |
|----|----|----|----|----|----|----|----|----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|



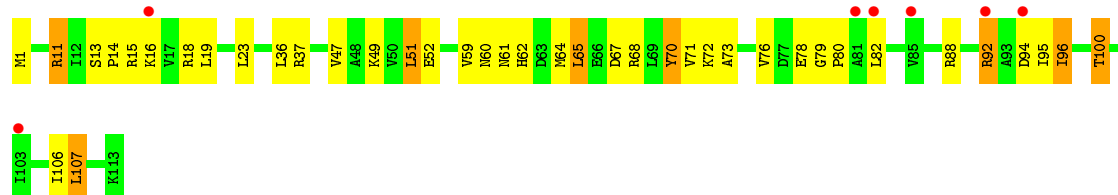
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|----|----|----|----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|
| M1 | T7 | G8 | G9 | K10 | Q11 | V14 | L18 | K19 | L20 | R21 | V22 | E23 | K24 | P29 | E34 | L35 | P36 | V37 | L38 | L39 | L40 | G41 | E43 | K44 | T45 | V46 | V47 | G48 | T49 | P50 | V51 | V52 | A55 | S56 | V57 | V58 | V61 | L62 | R66 | K69 | I70 | L71 | V72 | S73 | K74 | F75 | K76 | A77 | K78 |
|----|----|----|----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|



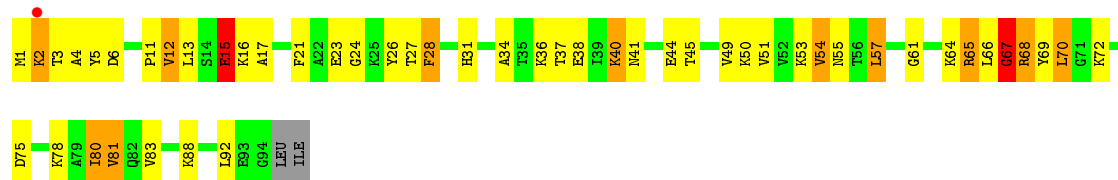
- Chain E8: 



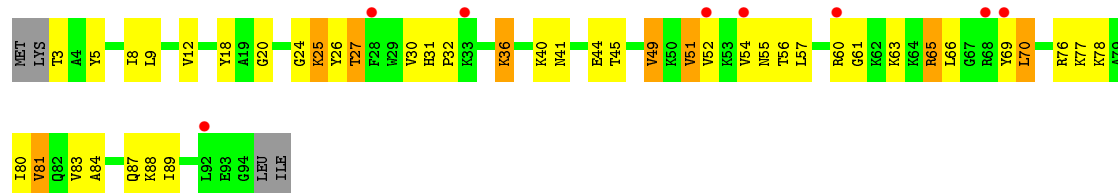
- Molecule 43: 50S ribosomal protein L22



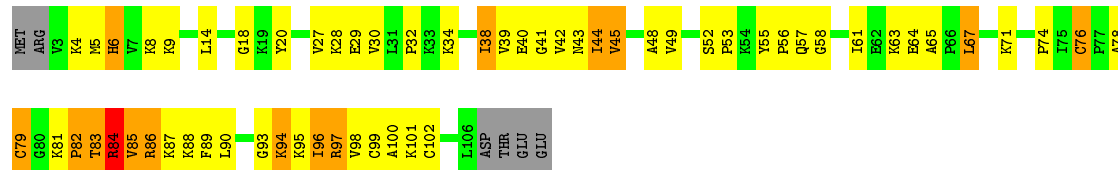
- Molecule 44: 50S ribosomal protein L23



- Molecule 44: 50S ribosomal protein L23

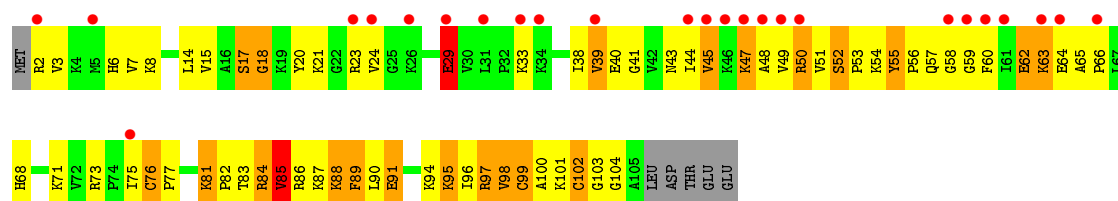


- Molecule 45: 50S ribosomal protein L24



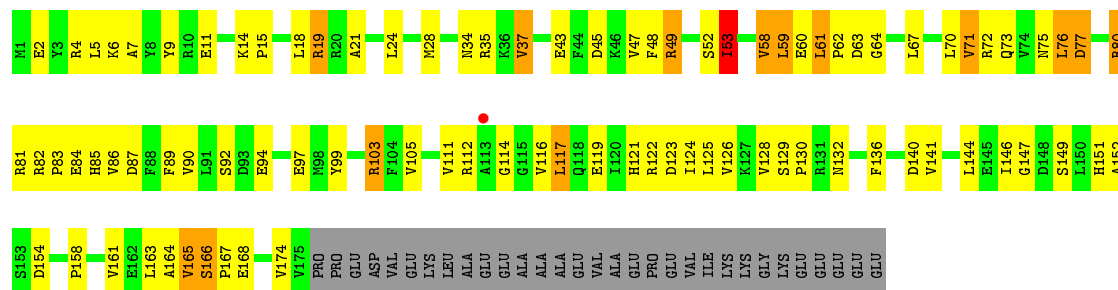
- Molecule 45: 50S ribosomal protein L24





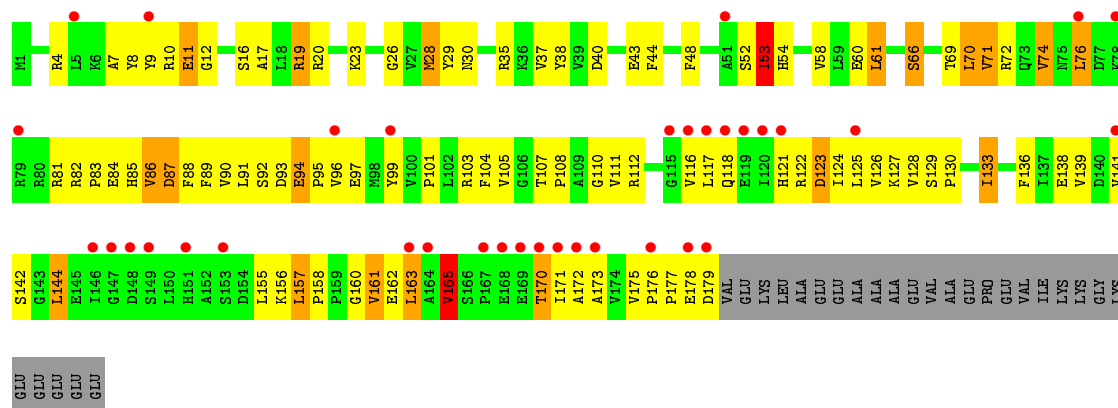
• Molecule 46: 50S ribosomal protein L25

Chain H8: 41% 36% 7% 15%



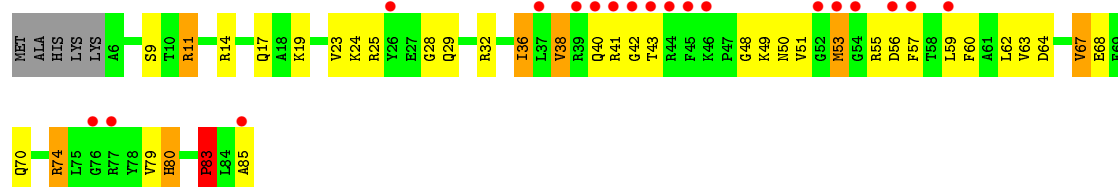
• Molecule 46: 50S ribosomal protein L25

Chain D5: 17% 38% 39% 9% 13%



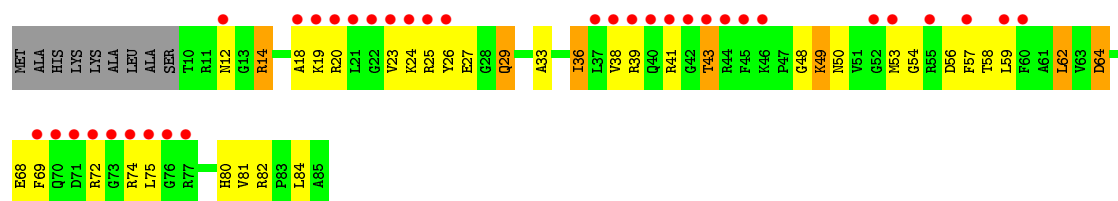
• Molecule 47: 50S ribosomal protein L27

Chain I8: 22% 49% 35% 8% 6%

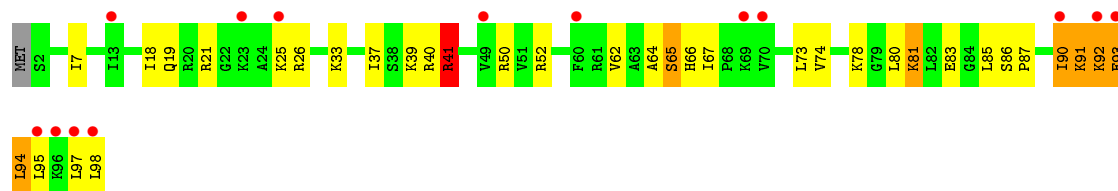


• Molecule 47: 50S ribosomal protein L27

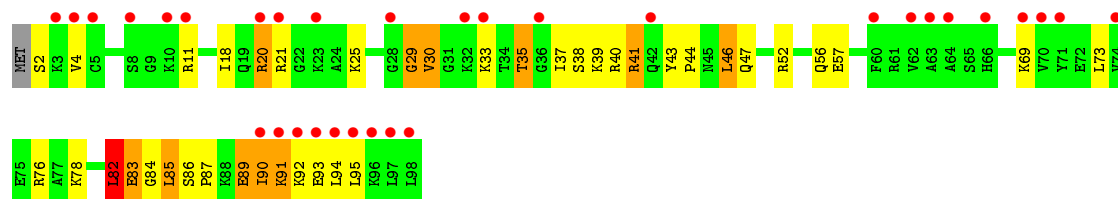
Chain E5: 41% 46% 35% 8% 11%



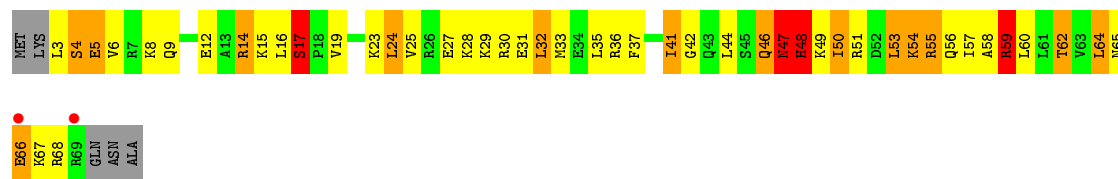
- Molecule 48: 50S ribosomal protein L28



- Molecule 48: 50S ribosomal protein L28



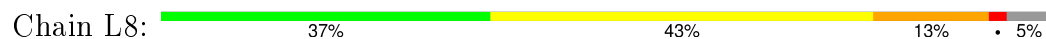
- Molecule 49: 50S ribosomal protein L29



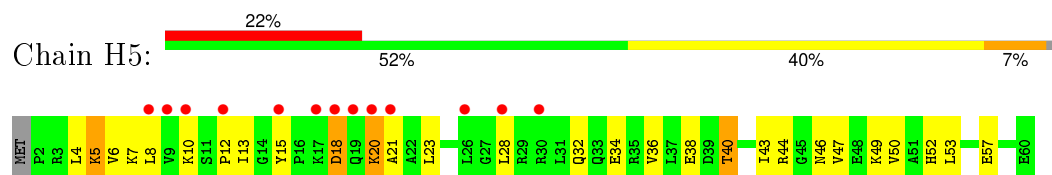
- Molecule 49: 50S ribosomal protein L29



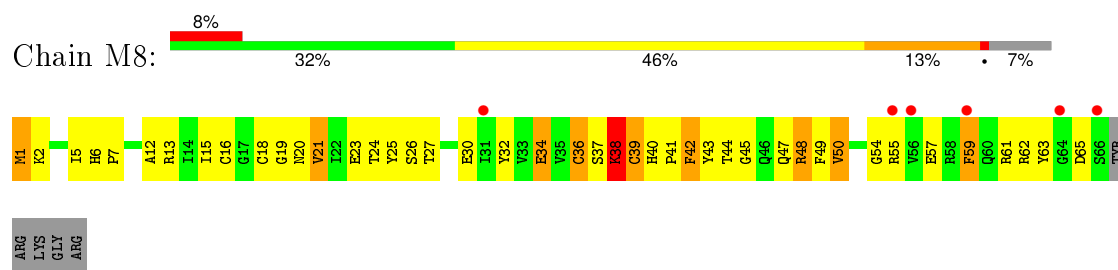
- Molecule 50: 50S ribosomal protein L30



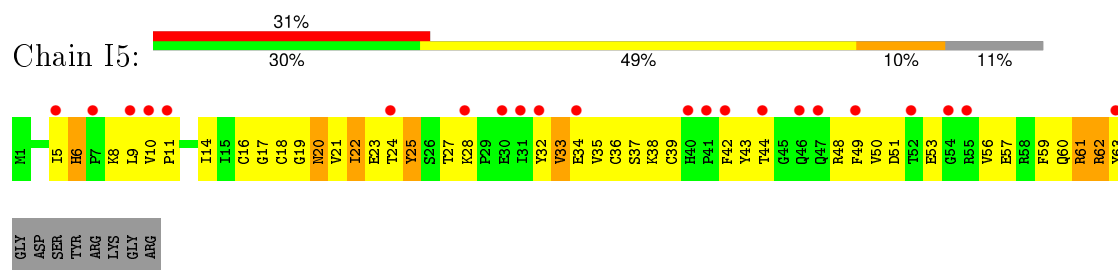
- Molecule 50: 50S ribosomal protein L30



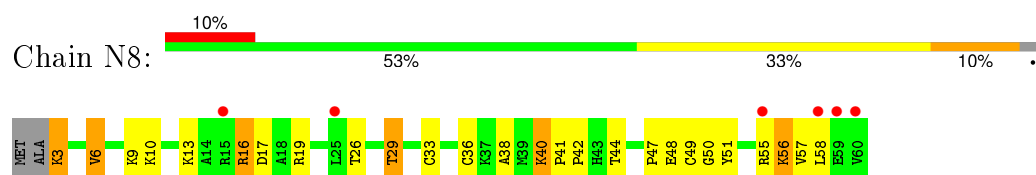
- Molecule 51: 50S ribosomal protein L31



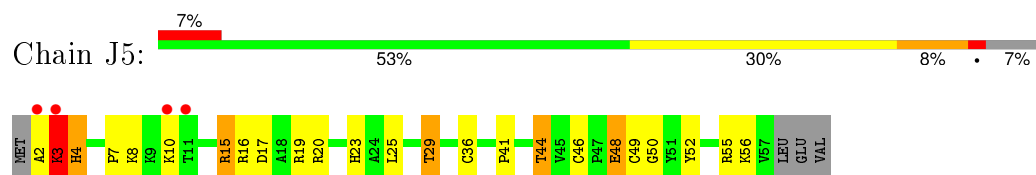
- Molecule 51: 50S ribosomal protein L31



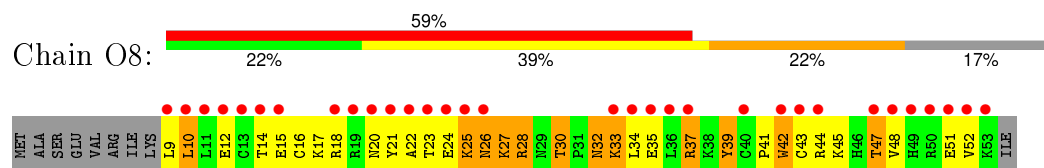
- Molecule 52: 50S ribosomal protein L32



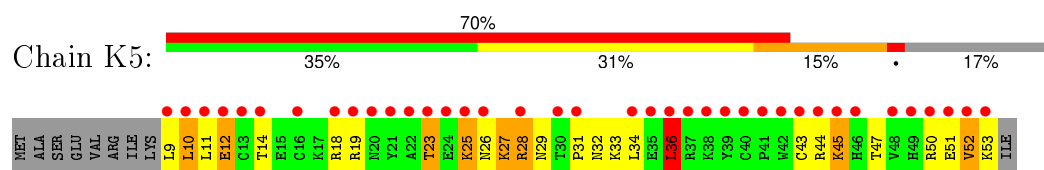
- Molecule 52: 50S ribosomal protein L32



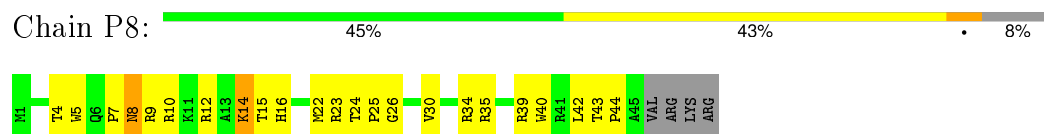
- Molecule 53: 50S ribosomal protein L33



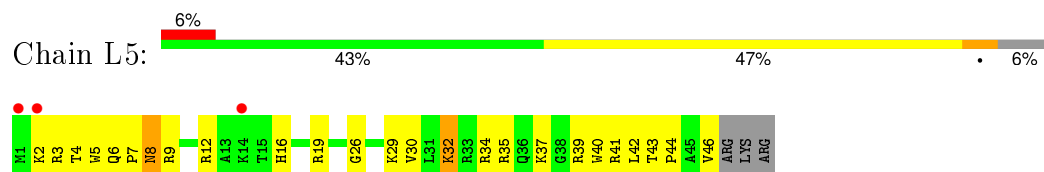
- Molecule 53: 50S ribosomal protein L33



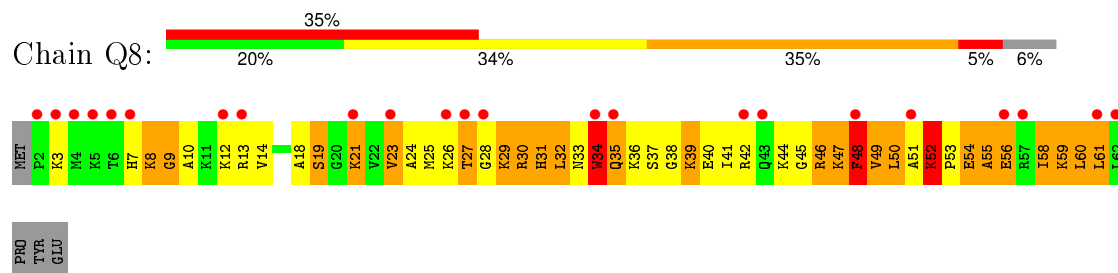
- Molecule 54: 50S ribosomal protein L34



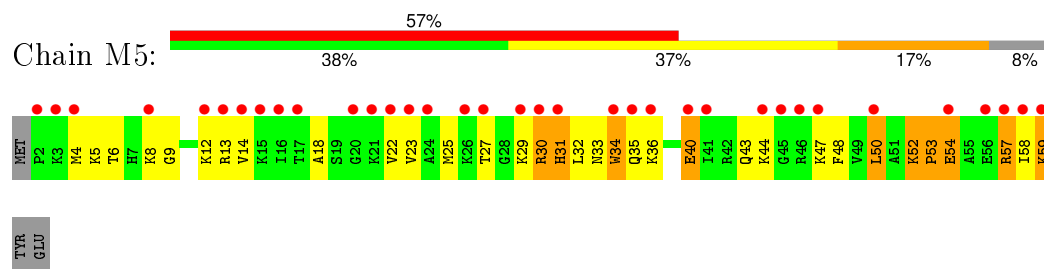
- Molecule 54: 50S ribosomal protein L34



- Molecule 55: 50S ribosomal protein L35



- Molecule 55: 50S ribosomal protein L35



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	209.90Å 447.90Å 621.50Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	188.03 – 3.05 223.95 – 3.05	Depositor EDS
% Data completeness (in resolution range)	99.8 (188.03-3.05) 91.4 (223.95-3.05)	Depositor EDS
R_{merge}	0.28	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.64 (at 3.07Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.199 , (Not available) 0.201 , 0.238	Depositor DCC
R_{free} test set	1998 reflections (0.20%)	DCC
Wilson B-factor (Å ²)	96.3	Xtriage
Anisotropy	0.265	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 89.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.27$	Xtriage
Outliers	0 of 1097157 reflections	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	299951	wwPDB-VP
Average B, all atoms (Å ²)	128.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.50% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: 5MU, OMC, ZN, MIA, MG, 4SU, 7MG, PSU

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	13	0.77	9/36053 (0.0%)	1.46	483/56270 (0.9%)
1	1G	0.65	1/36025 (0.0%)	1.30	249/56227 (0.4%)
2	12	0.37	0/1959	0.58	1/2642 (0.0%)
2	1E	0.42	0/1959	0.65	0/2642
3	22	0.41	0/1636	0.60	0/2205
3	2E	0.52	0/1629	0.67	0/2195
4	32	0.50	0/1732	0.73	1/2318 (0.0%)
4	3E	0.65	2/1732 (0.1%)	0.76	1/2318 (0.0%)
5	42	0.47	0/1171	0.74	0/1576
5	4E	0.55	0/1171	0.76	1/1576 (0.1%)
6	52	0.48	0/855	0.65	0/1154
6	5E	0.58	0/855	0.70	0/1154
7	62	0.42	0/1275	0.60	0/1709
7	6E	0.46	0/1275	0.59	0/1709
8	72	0.39	0/1135	0.60	0/1527
8	7E	0.51	0/1135	0.72	0/1527
9	82	0.39	0/1028	0.59	0/1379
9	8E	0.48	0/1028	0.72	1/1379 (0.1%)
10	1A	0.38	0/814	0.60	0/1095
10	1I	0.46	0/814	0.66	0/1095
11	2A	0.46	0/879	0.65	0/1187
11	2I	0.51	0/879	0.78	0/1187
12	3A	0.54	0/991	0.76	0/1327
12	3I	0.71	0/991	0.92	0/1327
13	4A	0.36	0/943	0.59	0/1265
13	4I	0.51	0/948	0.69	0/1272
14	5A	0.43	0/484	0.65	0/643
14	5I	0.71	1/500 (0.2%)	0.79	0/664
15	6A	0.47	0/744	0.62	1/992 (0.1%)
15	6I	0.54	0/744	0.78	0/992
16	7A	0.47	0/721	0.69	0/970
16	7I	0.48	0/721	0.72	0/970

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
17	8A	0.52	1/847 (0.1%)	0.64	0/1131
17	8I	0.57	0/847	0.74	0/1131
18	9A	0.51	0/595	0.77	0/790
18	9I	0.56	0/595	0.78	0/790
19	AA	0.40	0/638	0.63	0/860
19	AI	0.58	0/661	0.79	0/890
20	BA	0.48	0/764	0.76	0/1007
20	BI	0.45	0/764	0.70	0/1007
21	1B	0.44	0/221	0.64	0/288
21	1F	0.50	0/221	0.73	0/288
22	1K	0.51	0/1647	1.15	10/2565 (0.4%)
23	2K	1.12	6/1721 (0.3%)	1.64	40/2682 (1.5%)
23	2L	0.99	6/1721 (0.3%)	1.42	20/2682 (0.7%)
24	1L	0.44	0/1809	0.98	4/2819 (0.1%)
24	3K	0.49	1/1809 (0.1%)	1.17	14/2819 (0.5%)
24	3L	0.49	0/1809	1.11	11/2819 (0.4%)
25	4K	0.92	0/316	1.49	5/490 (1.0%)
25	4L	0.80	0/215	1.52	3/330 (0.9%)
26	14	0.91	64/70167 (0.1%)	1.65	1671/109541 (1.5%)
26	1H	1.09	145/70233 (0.2%)	1.89	2868/109643 (2.6%)
27	16	0.92	3/2928 (0.1%)	1.67	74/4568 (1.6%)
27	1J	0.75	0/2928	1.36	17/4568 (0.4%)
28	11	0.84	1/2165 (0.0%)	1.04	7/2919 (0.2%)
28	19	0.71	0/2170	0.96	4/2926 (0.1%)
29	21	0.68	0/1601	0.91	0/2160
29	29	0.69	0/1601	0.97	2/2160 (0.1%)
30	31	0.77	0/1620	0.97	2/2194 (0.1%)
30	39	0.61	0/1662	0.84	1/2249 (0.0%)
31	41	0.58	0/1498	0.83	1/2016 (0.0%)
31	49	0.41	0/1498	0.63	0/2016
32	51	0.66	0/1362	0.87	2/1841 (0.1%)
32	59	0.37	0/1324	0.63	0/1791
33	61	0.56	0/1151	0.79	2/1558 (0.1%)
33	69	0.47	0/1151	0.73	2/1558 (0.1%)
34	15	0.62	1/1131 (0.1%)	0.86	5/1525 (0.3%)
34	58	0.65	0/1131	0.86	3/1525 (0.2%)
35	25	0.64	0/942	0.77	0/1269
35	68	0.71	0/942	0.84	1/1269 (0.1%)
36	35	0.68	0/1161	1.06	4/1544 (0.3%)
36	78	0.75	0/1161	1.08	1/1544 (0.1%)
37	45	0.66	1/1142 (0.1%)	0.88	1/1527 (0.1%)
37	88	0.87	0/1106	1.05	2/1478 (0.1%)
38	55	0.66	0/973	0.88	2/1302 (0.2%)

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
38	98	0.60	0/981	0.87	0/1312
39	65	0.51	0/891	0.91	2/1187 (0.2%)
39	A8	0.71	0/891	0.98	4/1187 (0.3%)
40	75	0.59	0/1155	0.77	0/1542
40	B8	0.66	0/1155	0.84	1/1542 (0.1%)
41	85	0.57	0/981	0.75	1/1306 (0.1%)
41	C8	0.72	0/981	0.93	2/1306 (0.2%)
42	95	0.64	0/789	0.87	1/1057 (0.1%)
42	D8	0.66	0/789	0.84	1/1057 (0.1%)
43	A5	0.69	0/910	0.85	0/1220
43	E8	0.71	0/910	0.95	3/1220 (0.2%)
44	B5	0.77	0/739	0.86	0/993
44	F8	0.90	2/756 (0.3%)	0.98	0/1014
45	C5	0.69	0/807	0.90	1/1076 (0.1%)
45	G8	0.73	0/804	0.96	0/1073
46	D5	0.44	0/1460	0.67	0/1982
46	H8	0.50	0/1427	0.80	2/1935 (0.1%)
47	E5	0.66	0/614	0.88	0/819
47	I8	0.82	0/634	0.98	0/847
48	F5	0.66	0/769	0.93	2/1022 (0.2%)
48	J8	0.74	0/769	0.95	1/1022 (0.1%)
49	G5	0.58	0/560	0.81	1/741 (0.1%)
49	K8	0.94	2/565 (0.4%)	0.94	1/748 (0.1%)
50	H5	0.52	0/473	0.71	0/635
50	L8	0.73	1/457 (0.2%)	0.91	0/613
51	I5	0.46	0/527	0.72	0/709
51	M8	0.53	0/545	0.77	0/733
52	J5	0.71	0/448	0.88	0/606
52	N8	0.68	0/467	0.90	0/632
53	K5	0.74	0/396	0.85	1/529 (0.2%)
53	O8	0.87	1/396 (0.3%)	0.87	1/529 (0.2%)
54	L5	0.79	0/406	0.93	1/536 (0.2%)
54	P8	0.96	0/399	1.13	1/526 (0.2%)
55	M5	0.84	0/483	0.98	0/634
55	Q8	1.25	4/491 (0.8%)	1.62	7/645 (1.1%)
All	All	0.83	252/322599 (0.1%)	1.47	5551/483107 (1.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
4	3E	0	2
9	8E	0	1
10	1A	0	1
11	2I	0	2
13	4I	0	2
15	6I	0	1
16	7I	0	1
19	AI	0	2
20	BA	0	2
28	11	0	4
28	19	0	5
29	21	0	3
29	29	0	3
30	39	0	2
31	41	0	1
33	61	0	4
33	69	0	1
35	25	0	1
36	35	0	5
37	45	0	4
37	88	0	2
39	65	0	1
39	A8	0	2
40	75	0	2
40	B8	0	1
41	85	0	1
41	C8	0	1
43	E8	0	1
44	B5	0	2
44	F8	0	1
45	C5	0	3
45	G8	0	1
46	D5	0	1
46	H8	0	1
47	I8	0	2
48	F5	0	2
49	G5	0	3
49	K8	0	3
51	M8	0	2
52	J5	0	1
52	N8	0	1
53	O8	0	1
55	M5	0	1

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Mol	Chain	#Chirality outliers	#Planarity outliers
55	Q8	0	5
All	All	0	88

All (252) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
23	2L	21	U	C5-C6	19.03	1.51	1.34
23	2K	21	U	C5-C6	16.80	1.49	1.34
26	1H	774	A	N9-C4	-13.91	1.29	1.37
26	14	783	A	N9-C4	-12.95	1.30	1.37
23	2K	21	U	C2-N3	12.76	1.46	1.37
23	2L	21	U	C2-N3	12.46	1.46	1.37
26	1H	2430	A	N9-C4	-12.19	1.30	1.37
26	1H	783	A	N3-C4	-11.22	1.28	1.34
26	1H	676	A	N9-C4	-11.21	1.31	1.37
34	15	42	TRP	NE1-CE2	-10.77	1.23	1.37
26	1H	783	A	N9-C4	-10.71	1.31	1.37
26	1H	1614	A	N9-C4	-10.49	1.31	1.37
23	2K	21	U	N1-C2	10.36	1.47	1.38
26	1H	783	A	C5-C6	-10.32	1.31	1.41
23	2L	21	U	C4-C5	9.83	1.52	1.43
4	3E	12	CYS	CB-SG	9.57	1.98	1.82
26	1H	676	A	C5-C4	9.48	1.45	1.38
26	14	774	A	N9-C4	-9.41	1.32	1.37
26	1H	676	A	N9-C8	9.39	1.45	1.37
26	1H	621	A	N9-C4	-9.21	1.32	1.37
14	5I	27	CYS	CB-SG	-9.15	1.66	1.82
26	14	1950	G	C2-N3	9.05	1.40	1.32
26	1H	71	A	N9-C4	-8.97	1.32	1.37
26	1H	783	A	N7-C5	-8.94	1.33	1.39
23	2K	21	U	C4-C5	8.88	1.51	1.43
26	1H	1899	G	N9-C4	-8.75	1.30	1.38
26	1H	693	C	N3-C4	-8.58	1.27	1.33
23	2L	21	U	N1-C2	8.58	1.46	1.38
26	1H	1616	A	N9-C4	-8.56	1.32	1.37
26	1H	1786	A	C5-C6	-8.42	1.33	1.41
26	1H	805	G	N9-C8	-8.40	1.31	1.37
26	14	783	A	N3-C4	-8.40	1.29	1.34
26	1H	2346	A	N3-C4	-8.38	1.29	1.34
26	1H	1786	A	N7-C5	-8.34	1.34	1.39
49	K8	5	GLU	CG-CD	8.17	1.64	1.51
26	1H	1698	A	N9-C4	-8.15	1.32	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
26	14	1612	C	N1-C6	-8.13	1.32	1.37
26	1H	74	A	N9-C4	-7.92	1.33	1.37
26	14	783	A	C5-C6	-7.74	1.34	1.41
26	1H	1899	G	N3-C4	-7.71	1.30	1.35
23	2K	21	U	N3-C4	7.62	1.45	1.38
26	1H	945	A	C2-N3	7.53	1.40	1.33
26	1H	71	A	C6-N6	-7.52	1.27	1.33
26	14	945	A	C5-C6	-7.48	1.34	1.41
26	14	1616	A	N9-C4	-7.44	1.33	1.37
26	1H	1332	G	N9-C4	-7.42	1.32	1.38
1	13	792	A	N9-C4	-7.38	1.33	1.37
26	1H	2713	A	C5-C4	7.30	1.43	1.38
26	1H	676	A	N3-C4	-7.28	1.30	1.34
49	K8	5	GLU	CB-CG	7.27	1.66	1.52
53	O8	42	TRP	CB-CG	7.13	1.63	1.50
26	1H	1786	A	N3-C4	-6.98	1.30	1.34
26	1H	774	A	N3-C4	-6.97	1.30	1.34
23	2L	21	U	N3-C4	6.97	1.44	1.38
26	14	2346	A	N3-C4	-6.96	1.30	1.34
26	1H	2248	C	N3-C4	-6.94	1.29	1.33
23	2L	21	U	N1-C6	6.92	1.44	1.38
26	1H	860	U	N1-C2	6.92	1.44	1.38
26	1H	1786	A	N9-C4	-6.88	1.33	1.37
26	1H	1379	A	N9-C4	-6.87	1.33	1.37
26	14	74	A	N9-C4	-6.85	1.33	1.37
26	1H	945	A	N7-C5	-6.81	1.35	1.39
26	1H	1969	A	C6-N1	-6.80	1.30	1.35
55	Q8	54	GLU	CG-CD	6.79	1.62	1.51
26	1H	2252	G	C5-C4	-6.78	1.33	1.38
26	1H	828	U	C2-O2	6.75	1.28	1.22
26	1H	71	A	C5-C4	6.71	1.43	1.38
26	1H	138	G	N9-C8	6.70	1.42	1.37
26	1H	2585	U	N1-C2	6.66	1.44	1.38
26	1H	774	A	N9-C8	6.65	1.43	1.37
26	14	751	A	N9-C4	-6.64	1.33	1.37
26	14	676	A	N9-C4	-6.63	1.33	1.37
26	1H	1966	A	N9-C4	-6.62	1.33	1.37
4	3E	9	CYS	CB-SG	6.61	1.93	1.82
26	1H	945	A	N3-C4	-6.59	1.30	1.34
26	14	1272	A	N3-C4	6.58	1.38	1.34
26	1H	1899	G	N9-C8	6.58	1.42	1.37
26	1H	945	A	C5-C4	6.55	1.43	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
26	14	1950	G	C5-C4	6.54	1.43	1.38
26	1H	579	G	C2-N3	-6.51	1.27	1.32
23	2K	21	U	N1-C6	6.49	1.43	1.38
26	14	783	A	N7-C5	-6.49	1.35	1.39
1	13	792	A	N7-C5	-6.48	1.35	1.39
26	14	2287	A	N9-C4	-6.47	1.33	1.37
26	1H	71	A	N9-C8	6.46	1.43	1.37
26	1H	265	A	N3-C4	-6.42	1.31	1.34
26	1H	2031	A	N9-C4	6.41	1.41	1.37
26	14	2506	U	C2-N3	6.37	1.42	1.37
26	1H	74	A	N3-C4	-6.34	1.31	1.34
26	1H	528	A	N9-C4	-6.33	1.34	1.37
26	1H	1786	A	C5-C4	6.31	1.43	1.38
26	1H	1254	A	C5-C6	-6.30	1.35	1.41
27	16	115	G	C2-N3	6.27	1.37	1.32
26	1H	751	A	N3-C4	-6.27	1.31	1.34
26	1H	2287	A	N9-C4	-6.27	1.34	1.37
1	13	792	A	C5-C6	-6.25	1.35	1.41
26	14	1678	G	N9-C4	-6.24	1.32	1.38
26	14	2506	U	N1-C2	6.24	1.44	1.38
1	13	1498	U	N1-C2	6.22	1.44	1.38
26	1H	1614	A	N3-C4	-6.21	1.31	1.34
27	16	119	A	N9-C4	6.19	1.41	1.37
26	1H	829	A	N9-C4	-6.17	1.34	1.37
26	1H	2246	G	N9-C8	-6.12	1.33	1.37
26	14	1204	A	N9-C4	-6.10	1.34	1.37
26	1H	71	A	C5-C6	-6.09	1.35	1.41
26	1H	689	A	N3-C4	-6.08	1.31	1.34
26	1H	777	A	N3-C4	-6.08	1.31	1.34
28	11	122	ASP	CB-CG	6.05	1.64	1.51
26	1H	1142(A)	A	N9-C4	-6.04	1.34	1.37
55	Q8	56	GLU	CG-CD	6.03	1.60	1.51
26	14	1605	C	N1-C6	-6.02	1.33	1.37
26	1H	1621	U	N1-C6	-6.00	1.32	1.38
26	1H	2250	G	C2-N3	-6.00	1.27	1.32
1	13	1502	A	C5-C6	-5.99	1.35	1.41
26	1H	783	A	N9-C8	5.98	1.42	1.37
26	1H	1678	G	C5-C4	5.95	1.42	1.38
26	1H	805	G	N7-C5	-5.94	1.35	1.39
26	1H	192	C	N3-C4	5.94	1.38	1.33
26	1H	679	C	N1-C6	-5.92	1.33	1.37
26	1H	945	A	N9-C4	-5.92	1.34	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
26	1H	2594	C	N1-C6	-5.91	1.33	1.37
26	1H	1161	C	N1-C6	5.90	1.40	1.37
26	1H	216	A	N9-C4	-5.90	1.34	1.37
26	14	1617	C	N1-C6	-5.89	1.33	1.37
26	14	1819	A	N7-C5	-5.89	1.35	1.39
26	14	1303	G	C6-N1	-5.88	1.35	1.39
26	1H	1616	A	N7-C5	-5.88	1.35	1.39
26	1H	2373	G	C2-N3	5.87	1.37	1.32
26	1H	1815	A	N3-C4	-5.86	1.31	1.34
26	14	774	A	N9-C8	5.85	1.42	1.37
26	1H	2062	A	N3-C4	5.84	1.38	1.34
26	14	821	A	N7-C5	-5.83	1.35	1.39
26	1H	2071	A	C6-N6	-5.80	1.29	1.33
26	14	2518	A	N9-C4	-5.79	1.34	1.37
26	1H	746	A	N3-C4	-5.77	1.31	1.34
1	13	760	G	N7-C5	-5.77	1.35	1.39
26	14	528	A	N9-C4	-5.74	1.34	1.37
26	1H	138	G	C5-C4	5.74	1.42	1.38
26	14	1613	G	C8-N7	-5.74	1.27	1.30
50	L8	38	GLU	CG-CD	5.74	1.60	1.51
26	14	2447	G	N9-C8	-5.71	1.33	1.37
26	1H	1950	G	C2-N3	5.70	1.37	1.32
26	14	1786	A	C5-C6	-5.70	1.35	1.41
26	1H	2346	A	N9-C4	-5.70	1.34	1.37
24	3K	76	A	C5-C4	5.69	1.42	1.38
44	F8	15	GLU	CG-CD	5.68	1.60	1.51
26	14	465	G	N9-C4	5.68	1.42	1.38
26	1H	463	G	N1-C2	-5.68	1.33	1.37
26	14	216	A	N9-C4	-5.67	1.34	1.37
26	1H	245	G	N7-C5	-5.67	1.35	1.39
26	1H	789	A	N9-C4	-5.67	1.34	1.37
26	1H	2346	A	N7-C5	-5.67	1.35	1.39
44	F8	15	GLU	CB-CG	5.67	1.62	1.52
26	14	1950	G	N9-C4	5.66	1.42	1.38
26	1H	1678	G	N9-C8	5.66	1.41	1.37
26	1H	621	A	C5-C6	-5.65	1.35	1.41
26	1H	821	A	N7-C5	-5.64	1.35	1.39
26	14	676	A	N9-C8	5.64	1.42	1.37
26	14	1950	G	N3-C4	5.64	1.39	1.35
26	1H	213	A	N9-C4	-5.63	1.34	1.37
26	1H	1692	U	C4-C5	-5.63	1.38	1.43
26	1H	774	A	C8-N7	5.63	1.35	1.31

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
26	1H	2254	C	N1-C2	-5.61	1.34	1.40
26	1H	1332	G	C5-C4	5.59	1.42	1.38
26	14	2332	U	C4-C5	5.58	1.48	1.43
26	1H	195	A	C6-N1	5.57	1.39	1.35
26	1H	203	C	N1-C2	-5.56	1.34	1.40
26	14	1396	U	N1-C2	5.55	1.43	1.38
26	1H	945	A	N1-C2	5.54	1.39	1.34
26	1H	52	A	N9-C4	5.54	1.41	1.37
26	14	1278	A	N9-C4	-5.54	1.34	1.37
26	14	2346	A	N9-C4	-5.54	1.34	1.37
27	16	101	A	N9-C4	-5.53	1.34	1.37
26	1H	1427	A	N9-C4	5.51	1.41	1.37
37	45	80	GLU	CG-CD	5.51	1.60	1.51
26	1H	1382	G	C5-C6	-5.49	1.36	1.42
26	14	2606	C	N1-C6	-5.49	1.33	1.37
26	1H	204	A	N3-C4	-5.49	1.31	1.34
26	1H	1255	U	C2-N3	5.47	1.41	1.37
26	1H	70	G	C6-N1	-5.46	1.35	1.39
26	14	2430	A	N9-C4	-5.46	1.34	1.37
26	1H	2440	C	N1-C6	-5.44	1.33	1.37
26	1H	451	C	N1-C6	-5.43	1.33	1.37
26	1H	2256	G	N7-C5	-5.42	1.35	1.39
26	14	1997	G	C2-N3	5.42	1.37	1.32
26	1H	1255	U	C4-O4	5.41	1.27	1.23
26	14	1332	G	C5-C4	5.41	1.42	1.38
26	14	2518	A	C5-C6	-5.40	1.36	1.41
26	14	1950	G	C5-C6	5.40	1.47	1.42
1	1G	1473	A	N9-C4	-5.38	1.34	1.37
26	1H	1957	C	N3-C4	-5.38	1.30	1.33
26	1H	2502	G	N1-C2	-5.38	1.33	1.37
26	14	1676	A	N3-C4	-5.36	1.31	1.34
26	14	472	A	N3-C4	-5.36	1.31	1.34
26	14	1660	C	N3-C4	-5.33	1.30	1.33
26	14	2249	U	C4-O4	5.33	1.27	1.23
26	1H	140	A	C5-C6	-5.32	1.36	1.41
26	1H	608	A	N3-C4	-5.32	1.31	1.34
26	14	1698	A	C5-C6	-5.32	1.36	1.41
26	14	2062	A	C6-N1	5.32	1.39	1.35
26	14	775	G	C6-O6	-5.31	1.19	1.24
55	Q8	48	PHE	CB-CG	-5.30	1.42	1.51
26	1H	783	A	C6-N1	-5.29	1.31	1.35
26	1H	2377	A	N9-C4	-5.29	1.34	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
26	1H	2062	A	N7-C5	5.28	1.42	1.39
26	1H	2713	A	N9-C4	-5.28	1.34	1.37
26	14	204	A	N7-C5	-5.28	1.36	1.39
26	1H	774	A	C2-N3	-5.27	1.28	1.33
26	1H	2602	A	N3-C4	5.27	1.38	1.34
26	1H	2430	A	C5-C6	-5.27	1.36	1.41
26	1H	1210	A	N9-C4	-5.26	1.34	1.37
26	14	2612	C	N3-C4	5.26	1.37	1.33
26	1H	470	A	N3-C4	-5.26	1.31	1.34
26	1H	698	C	N1-C6	-5.25	1.34	1.37
26	1H	1613	G	N7-C5	-5.24	1.36	1.39
1	13	690	G	C5-C4	5.24	1.42	1.38
26	1H	692	C	C2-N3	5.24	1.40	1.35
26	1H	140	A	N7-C5	-5.23	1.36	1.39
26	1H	777	A	N9-C4	-5.23	1.34	1.37
26	14	2015	A	N3-C4	-5.23	1.31	1.34
26	1H	1982	C	C4-C5	-5.23	1.38	1.43
26	1H	2713	A	N9-C8	5.23	1.42	1.37
26	1H	1784	A	N7-C5	5.23	1.42	1.39
26	1H	1392	A	N9-C4	5.22	1.41	1.37
26	14	1645	G	C5-C4	-5.21	1.34	1.38
26	14	676	A	C5-C4	5.19	1.42	1.38
26	1H	1698	A	C5-C6	-5.18	1.36	1.41
26	1H	2053	G	C5-C4	-5.18	1.34	1.38
26	1H	1888	G	N9-C4	5.17	1.42	1.38
26	1H	2494	G	C2-N3	-5.16	1.28	1.32
26	14	1786	A	N9-C4	-5.14	1.34	1.37
26	1H	187	G	N7-C5	-5.14	1.36	1.39
26	1H	1311	G	N9-C8	-5.14	1.34	1.37
26	1H	2451	A	C6-N1	-5.14	1.31	1.35
26	14	621	A	N9-C4	-5.13	1.34	1.37
26	1H	952	G	C5-C4	-5.12	1.34	1.38
26	14	1825	A	C6-N1	-5.12	1.31	1.35
26	14	1698	A	N7-C5	-5.12	1.36	1.39
1	13	1227	A	N9-C4	-5.11	1.34	1.37
26	1H	1385	G	N9-C4	-5.11	1.33	1.38
26	1H	1971	A	C5-C4	-5.11	1.35	1.38
55	Q8	54	GLU	CB-CG	5.11	1.61	1.52
26	1H	1786	A	N1-C2	5.11	1.39	1.34
26	1H	122	G	C2-N3	5.10	1.36	1.32
1	13	1483	A	N9-C4	-5.09	1.34	1.37
26	1H	1786	A	C6-N1	-5.08	1.31	1.35

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
26	1H	1950	G	N9-C8	5.08	1.41	1.37
26	14	515	A	N9-C4	-5.08	1.34	1.37
26	1H	663	G	N3-C4	-5.05	1.31	1.35
17	8A	49	GLU	CG-CD	5.04	1.59	1.51
26	14	1342	A	N3-C4	-5.04	1.31	1.34
26	1H	1349	A	N9-C8	5.03	1.41	1.37

All (5551) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	1786	A	C2-N3-C4	-22.74	99.23	110.60
26	1H	1899	G	N3-C4-N9	-22.01	112.80	126.00
26	1H	676	A	C2-N3-C4	-20.28	100.46	110.60
26	1H	1899	G	C2-N3-C4	-19.38	102.21	111.90
26	1H	783	A	C2-N3-C4	-18.91	101.14	110.60
26	14	741	G	O5'-P-OP1	-18.63	88.34	110.70
26	1H	1786	A	N1-C2-N3	17.51	138.05	129.30
26	1H	1899	G	N3-C4-C5	17.49	137.35	128.60
26	1H	1982	C	O5'-P-OP2	-17.27	89.98	110.70
26	1H	1332	G	C2-N3-C4	-17.00	103.40	111.90
26	1H	1786	A	N7-C8-N9	16.89	122.25	113.80
26	1H	945	A	C4-C5-C6	16.82	125.41	117.00
26	1H	945	A	C6-C5-N7	-16.71	120.60	132.30
26	1H	774	A	N3-C4-N9	-16.44	114.25	127.40
26	1H	968	G	O5'-P-OP2	-16.32	91.01	105.70
26	1H	783	A	C5-N7-C8	-16.11	95.85	103.90
26	1H	74	A	C2-N3-C4	-15.94	102.63	110.60
26	14	1332	G	C6-C5-N7	-15.90	120.86	130.40
26	1H	1614	A	C2-N3-C4	-15.60	102.80	110.60
26	14	2436	G	O5'-P-OP1	-15.60	91.66	105.70
26	14	783	A	C2-N3-C4	-15.26	102.97	110.60
26	1H	1786	A	C5-N7-C8	-15.25	96.27	103.90
26	1H	2430	A	N1-C6-N6	15.12	127.67	118.60
26	14	945	A	N1-C6-N6	14.88	127.53	118.60
26	14	74	A	C2-N3-C4	-14.87	103.17	110.60
26	1H	1829	A	O5'-P-OP1	-14.86	92.33	105.70
26	1H	774	A	N3-C4-C5	14.76	137.13	126.80
26	1H	783	A	C8-N9-C4	-14.73	99.91	105.80
26	1H	676	A	C5-N7-C8	-14.63	96.58	103.90
26	1H	621	A	C5-N7-C8	-14.60	96.60	103.90
26	1H	621	A	C2-N3-C4	-14.57	103.31	110.60
26	1H	71	A	C2-N3-C4	-14.57	103.32	110.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	1786	A	C6-C5-N7	-14.55	122.12	132.30
26	1H	1678	G	C2-N3-C4	-14.42	104.69	111.90
26	1H	2430	A	O5'-P-OP2	-14.37	92.77	105.70
26	1H	917	A	C2-N3-C4	-14.31	103.44	110.60
26	1H	1779	U	O5'-P-OP2	-14.27	92.86	105.70
26	1H	2314	C	O5'-P-OP2	-14.23	92.89	105.70
26	14	1698	A	N1-C6-N6	14.12	127.07	118.60
26	1H	1678	G	C5-N7-C8	-14.08	97.26	104.30
26	1H	1332	G	C5-N7-C8	-14.03	97.29	104.30
26	14	783	A	N1-C6-N6	14.01	127.00	118.60
26	1H	676	A	N3-C4-N9	-13.91	116.27	127.40
26	1H	945	A	N7-C8-N9	13.91	120.75	113.80
26	14	1899	G	N1-C2-N2	-13.88	103.71	116.20
26	1H	2430	A	C2-N3-C4	-13.85	103.68	110.60
26	14	530	G	N1-C6-O6	13.65	128.09	119.90
26	1H	676	A	N7-C8-N9	13.64	120.62	113.80
26	1H	1312	U	O5'-P-OP1	-13.63	93.44	105.70
26	1H	860	U	N3-C2-O2	-13.55	112.72	122.20
26	14	2430	A	N1-C6-N6	13.54	126.73	118.60
26	1H	621	A	N1-C6-N6	13.51	126.70	118.60
26	1H	860	U	C4-C5-C6	13.47	127.78	119.70
26	1H	801	G	O5'-P-OP2	-13.44	93.61	105.70
26	1H	676	A	N3-C4-C5	13.42	136.19	126.80
26	1H	917	A	N1-C2-N3	13.37	135.98	129.30
26	1H	2287	A	C2-N3-C4	-13.35	103.92	110.60
26	1H	1786	A	C8-N9-C4	-13.34	100.46	105.80
26	1H	774	A	C5-N7-C8	-13.25	97.27	103.90
26	1H	774	A	C6-N1-C2	13.23	126.54	118.60
26	14	1786	A	C2-N3-C4	-13.21	104.00	110.60
26	1H	140	A	N7-C8-N9	13.18	120.39	113.80
26	14	530	G	N9-C4-C5	-13.15	100.14	105.40
26	14	2518	A	N1-C6-N6	13.14	126.48	118.60
26	1H	1678	G	N7-C8-N9	13.04	119.62	113.10
26	1H	1698	A	C2-N3-C4	-13.04	104.08	110.60
26	1H	2490	G	C5-N7-C8	-12.97	97.81	104.30
23	2K	61	U	O5'-P-OP2	-12.96	94.04	105.70
26	1H	733	G	O5'-P-OP2	-12.92	94.07	105.70
26	1H	1332	G	N3-C4-C5	12.91	135.06	128.60
26	1H	2346	A	N1-C2-N3	12.89	135.75	129.30
26	1H	1616	A	C5-N7-C8	-12.89	97.46	103.90
26	1H	1204	A	O4'-C1'-N9	12.86	118.49	108.20
26	14	1332	G	C4-C5-N7	12.80	115.92	110.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	783	A	N7-C8-N9	12.79	120.19	113.80
26	14	1639	U	O5'-P-OP2	-12.78	94.20	105.70
26	14	1274	A	N1-C6-N6	12.77	126.26	118.60
26	14	1786	A	C5-N7-C8	-12.76	97.52	103.90
26	1H	1931	U	N3-C2-O2	-12.71	113.31	122.20
26	1H	793	A	O5'-P-OP2	-12.69	94.28	105.70
26	1H	945	A	N1-C2-N3	12.68	135.64	129.30
26	14	676	A	C2-N3-C4	-12.67	104.27	110.60
23	2K	21	U	C5-C4-O4	-12.64	118.31	125.90
26	1H	120	U	C5-C6-N1	-12.63	116.38	122.70
26	1H	140	A	C5-N7-C8	-12.63	97.59	103.90
26	1H	51	G	O5'-P-OP1	-12.57	94.39	105.70
26	1H	1382	G	C5-C6-O6	-12.54	121.08	128.60
26	14	1786	A	N7-C8-N9	12.52	120.06	113.80
26	1H	2430	A	C5-N7-C8	-12.49	97.66	103.90
26	1H	2699	C	C6-N1-C2	12.44	125.28	120.30
26	1H	1772	G	N1-C6-O6	-12.44	112.44	119.90
26	14	774	A	N3-C4-N9	-12.36	117.51	127.40
26	1H	71	A	C5-N7-C8	-12.35	97.73	103.90
26	14	774	A	N3-C4-C5	12.34	135.44	126.80
26	14	1332	G	C5-N7-C8	-12.34	98.13	104.30
26	14	783	A	C5-N7-C8	-12.27	97.77	103.90
1	13	792	A	O4'-C1'-N9	12.25	118.00	108.20
26	1H	1899	G	N3-C2-N2	-12.24	111.33	119.90
26	1H	676	A	C5-C6-N1	-12.22	111.59	117.70
26	1H	1678	G	C4-C5-N7	12.20	115.68	110.80
26	1H	945	A	N1-C6-N6	12.15	125.89	118.60
26	1H	2346	A	O4'-C1'-N9	12.15	117.92	108.20
26	14	2430	A	C2-N3-C4	-12.15	104.53	110.60
26	1H	46	C	O5'-P-OP2	-12.14	94.78	105.70
26	1H	2713	A	N7-C8-N9	12.10	119.85	113.80
26	14	2873	A	N7-C8-N9	12.09	119.85	113.80
1	13	690	G	C6-C5-N7	-12.08	123.15	130.40
26	1H	1393	A	O5'-P-OP2	-12.04	94.87	105.70
26	14	1602	U	O5'-P-OP2	12.02	125.13	110.70
24	3K	76	A	C5-N7-C8	-12.00	97.90	103.90
26	14	1332	G	N1-C6-O6	11.96	127.08	119.90
26	1H	133	C	C6-N1-C2	11.94	125.08	120.30
26	1H	2053	G	C5-C6-O6	-11.94	121.44	128.60
26	14	1332	G	N7-C8-N9	11.94	119.07	113.10
26	14	1899	G	N3-C2-N2	11.89	128.22	119.90
24	3K	76	A	N7-C8-N9	11.88	119.74	113.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	140	A	C8-N9-C4	-11.87	101.05	105.80
26	1H	2713	A	C5-N7-C8	-11.87	97.97	103.90
26	1H	945	A	C5-N7-C8	-11.77	98.01	103.90
26	1H	783	A	C4-C5-N7	11.72	116.56	110.70
26	14	828	U	C5-C4-O4	11.61	132.86	125.90
26	14	676	A	C5-N7-C8	-11.60	98.10	103.90
26	14	530	G	C5-C6-O6	-11.58	121.65	128.60
26	1H	2490	G	C4-C5-N7	11.56	115.42	110.80
26	1H	783	A	C6-C5-N7	-11.55	124.21	132.30
1	13	690	G	C4-N9-C1'	11.54	141.50	126.50
26	14	945	A	C6-C5-N7	-11.53	124.23	132.30
26	1H	1647	G	O5'-P-OP1	-11.53	95.33	105.70
26	1H	621	A	C4-C5-N7	11.52	116.46	110.70
26	1H	200	U	O5'-P-OP1	-11.51	95.34	105.70
26	14	945	A	C2-N3-C4	-11.51	104.85	110.60
26	1H	676	A	C8-N9-C4	-11.49	101.20	105.80
26	1H	945	A	C2-N3-C4	-11.49	104.85	110.60
26	14	922	U	O5'-P-OP1	-11.48	95.37	105.70
26	1H	2430	A	N3-C4-C5	11.48	134.84	126.80
26	1H	391	G	N1-C6-O6	11.48	126.79	119.90
26	14	945	A	C4-C5-N7	11.46	116.43	110.70
26	1H	530	G	N1-C6-O6	-11.46	113.03	119.90
26	1H	860	U	C5-C6-N1	-11.43	116.99	122.70
26	1H	624	C	O5'-P-OP2	11.43	124.41	110.70
26	1H	2265	U	O5'-P-OP1	-11.39	95.45	105.70
26	1H	2710	C	C6-N1-C2	11.38	124.85	120.30
26	1H	913	U	N1-C2-O2	11.36	130.75	122.80
26	1H	966	G	O5'-P-OP2	-11.34	95.50	105.70
26	1H	1249	U	O5'-P-OP1	-11.32	95.51	105.70
26	14	1698	A	C6-C5-N7	-11.31	124.38	132.30
26	1H	1616	A	N7-C8-N9	11.30	119.45	113.80
24	3L	76	A	N7-C8-N9	11.23	119.42	113.80
26	1H	1604	C	N1-C2-O2	-11.20	112.18	118.90
26	14	530	G	C8-N9-C4	11.16	110.86	106.40
26	1H	120	U	C4-C5-C6	11.15	126.39	119.70
26	14	2873	A	C5-N7-C8	-11.15	98.33	103.90
26	1H	2598	A	O5'-P-OP1	-11.13	95.68	105.70
26	1H	1336	A	N1-C6-N6	-11.12	111.93	118.60
26	1H	1899	G	N9-C4-C5	11.12	109.85	105.40
26	1H	2070	G	N1-C2-N2	-11.07	106.23	116.20
26	1H	2380	C	C5-C6-N1	-11.02	115.49	121.00
26	14	2778	A	O5'-P-OP2	-11.01	95.79	105.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	1678	G	C6-C5-N7	-11.01	123.80	130.40
26	1H	446	G	N9-C4-C5	-11.01	101.00	105.40
1	13	792	A	C5-N7-C8	-11.00	98.40	103.90
26	1H	1496	A	C5-N7-C8	-11.00	98.40	103.90
26	1H	216	A	O5'-P-OP1	-10.98	95.82	105.70
26	14	2441	C	O5'-P-OP1	-10.93	95.86	105.70
26	1H	789	A	O5'-P-OP1	-10.93	95.87	105.70
26	1H	2070	G	N3-C2-N2	10.93	127.55	119.90
26	1H	827	U	O5'-P-OP2	-10.91	95.88	105.70
26	1H	799	G	C8-N9-C4	10.90	110.76	106.40
26	1H	2374	C	C5-C6-N1	-10.89	115.55	121.00
26	1H	1899	G	N1-C2-N3	10.89	130.43	123.90
26	1H	1614	A	N1-C2-N3	10.88	134.74	129.30
1	13	813	U	O5'-P-OP2	-10.87	95.92	105.70
26	1H	2062	A	C8-N9-C4	10.85	110.14	105.80
26	1H	2819	G	N1-C6-O6	10.85	126.41	119.90
26	14	1619	G	O5'-P-OP2	-10.84	95.94	105.70
26	1H	1786	A	N1-C6-N6	10.82	125.09	118.60
26	1H	1332	G	C4-C5-N7	10.82	115.13	110.80
26	1H	1332	G	N7-C8-N9	10.79	118.50	113.10
26	1H	613	U	N3-C2-O2	-10.78	114.65	122.20
26	1H	2518	A	C5-N7-C8	-10.74	98.53	103.90
26	1H	1787	A	O5'-P-OP1	-10.72	96.05	105.70
26	1H	2247	A	O5'-P-OP1	-10.72	96.05	105.70
26	1H	1496	A	N7-C8-N9	10.70	119.15	113.80
26	1H	2584	U	N3-C2-O2	-10.70	114.71	122.20
26	14	793	A	O5'-P-OP2	-10.70	96.07	105.70
26	14	1984	G	O5'-P-OP2	-10.69	96.08	105.70
26	1H	2295	C	C6-N1-C2	-10.68	116.03	120.30
26	14	2688	U	N3-C2-O2	-10.67	114.73	122.20
26	1H	945	A	C8-N9-C4	-10.67	101.53	105.80
23	2K	21	U	N3-C4-C5	10.63	120.98	114.60
26	14	1275	A	O5'-P-OP1	-10.60	96.16	105.70
26	14	2518	A	C5-N7-C8	-10.58	98.61	103.90
26	1H	2700	C	C6-N1-C2	10.57	124.53	120.30
26	14	312	G	O5'-P-OP1	-10.55	96.20	105.70
26	1H	1786	A	C4-C5-C6	10.54	122.27	117.00
26	1H	2507	C	N3-C2-O2	-10.54	114.53	121.90
26	1H	2606	C	O5'-P-OP1	-10.53	96.22	105.70
26	1H	2713	A	C8-N9-C4	-10.53	101.59	105.80
26	1H	774	A	C5-C6-N1	-10.52	112.44	117.70
26	1H	1379	A	N1-C6-N6	10.50	124.90	118.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	863	A	O5'-P-OP2	-10.49	96.26	105.70
26	1H	2698	U	O5'-P-OP2	-10.47	96.28	105.70
26	1H	1931	U	N1-C2-N3	10.46	121.18	114.90
26	1H	621	A	N7-C8-N9	10.45	119.02	113.80
26	14	945	A	C5-N7-C8	-10.45	98.68	103.90
26	14	687	C	O5'-P-OP1	-10.45	96.30	105.70
26	1H	783	A	N3-C4-N9	-10.42	119.06	127.40
26	14	689	A	O5'-P-OP2	-10.41	96.33	105.70
26	1H	2392	A	C5-N7-C8	-10.41	98.70	103.90
26	1H	265	A	C8-N9-C4	-10.39	101.64	105.80
26	1H	252	G	O5'-P-OP2	-10.39	96.35	105.70
26	1H	1566	A	O5'-P-OP2	-10.38	96.36	105.70
26	14	1816	G	O5'-P-OP1	-10.36	96.38	105.70
26	14	2702	U	O5'-P-OP2	-10.36	96.38	105.70
26	14	2430	A	C5-C6-N1	-10.35	112.52	117.70
27	16	47	C	C6-N1-C2	10.34	124.44	120.30
26	1H	330	A	C2-N3-C4	-10.34	105.43	110.60
26	14	929	G	N1-C6-O6	10.34	126.10	119.90
26	14	1616	A	C5-N7-C8	-10.33	98.73	103.90
26	1H	2036	C	C6-N1-C2	-10.33	116.17	120.30
26	1H	528	A	C6-N1-C2	10.32	124.79	118.60
26	1H	1437	C	C6-N1-C2	-10.29	116.19	120.30
26	1H	582	G	N1-C6-O6	10.26	126.06	119.90
26	14	963	U	O5'-P-OP1	-10.26	96.46	105.70
26	14	530	G	C4-C5-N7	10.26	114.90	110.80
26	1H	1528	A	C8-N9-C4	-10.25	101.70	105.80
26	14	1678	G	C5-N7-C8	-10.25	99.18	104.30
26	14	783	A	N3-C4-C5	10.24	133.97	126.80
26	1H	2311	A	C2-N3-C4	-10.24	105.48	110.60
1	13	789	U	C5-C4-O4	10.23	132.04	125.90
26	1H	463	G	N3-C2-N2	10.22	127.06	119.90
26	1H	461	C	N1-C2-O2	-10.22	112.77	118.90
26	14	140	A	C5-N7-C8	-10.20	98.80	103.90
26	1H	945	A	C4-N9-C1'	10.20	144.65	126.30
1	13	690	G	C8-N9-C1'	-10.19	113.75	127.00
26	14	2429	G	O5'-P-OP1	10.18	122.92	110.70
26	1H	802	A	O5'-P-OP2	-10.18	96.54	105.70
26	14	1332	G	C2-N3-C4	-10.18	106.81	111.90
26	1H	74	A	N1-C2-N3	10.18	134.39	129.30
26	1H	1616	A	C8-N9-C4	-10.16	101.74	105.80
26	1H	676	A	O4'-C1'-N9	10.15	116.32	108.20
26	14	2440	C	C6-N1-C2	10.13	124.35	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	1281	G	O5'-P-OP2	10.13	122.86	110.70
1	1G	1395	C	O5'-P-OP1	-10.13	96.58	105.70
26	1H	765	G	O5'-P-OP1	-10.12	96.59	105.70
26	1H	624	C	O5'-P-OP1	-10.11	96.60	105.70
26	1H	812	C	N1-C2-O2	-10.11	112.84	118.90
26	14	783	A	C4-C5-N7	10.10	115.75	110.70
1	13	580	U	C5-C6-N1	-10.09	117.65	122.70
26	14	528	A	C2-N3-C4	-10.08	105.56	110.60
26	1H	49	A	O5'-P-OP2	-10.08	96.63	105.70
26	1H	769	G	N3-C2-N2	10.06	126.94	119.90
26	14	1698	A	C2-N3-C4	-10.06	105.57	110.60
26	14	2610	C	C6-N1-C2	10.06	124.32	120.30
26	1H	195	A	N1-C6-N6	10.05	124.63	118.60
26	1H	1394	U	O5'-P-OP1	-10.05	96.65	105.70
1	13	899	C	N1-C2-O2	-10.04	112.88	118.90
26	14	1367	A	N1-C6-N6	10.04	124.62	118.60
26	14	2712	U	C5-C6-N1	-10.03	117.69	122.70
26	14	1253	A	N1-C6-N6	9.99	124.59	118.60
26	1H	2346	A	C2-N3-C4	-9.98	105.61	110.60
26	1H	1825	A	N1-C6-N6	-9.97	112.62	118.60
26	1H	2430	A	O5'-P-OP1	9.97	122.66	110.70
1	13	792	A	C4-C5-N7	9.95	115.67	110.70
26	14	2244	U	N1-C2-O2	-9.94	115.84	122.80
26	1H	2747	G	N1-C6-O6	9.93	125.86	119.90
26	14	2430	A	N1-C2-N3	9.93	134.26	129.30
26	14	74	A	N1-C6-N6	9.91	124.55	118.60
26	1H	678	C	C2-N3-C4	-9.91	114.95	119.90
26	1H	2490	G	N7-C8-N9	9.90	118.05	113.10
24	3L	76	A	C5-N7-C8	-9.90	98.95	103.90
26	1H	2618	G	C8-N9-C4	-9.90	102.44	106.40
26	1H	1950	G	C5-N7-C8	-9.87	99.36	104.30
26	1H	2430	A	C4-C5-N7	9.86	115.63	110.70
26	1H	265	A	C2-N3-C4	-9.84	105.68	110.60
26	1H	739	G	O5'-P-OP2	-9.83	96.86	105.70
26	1H	913	U	N3-C4-O4	-9.83	112.52	119.40
26	1H	1210	A	C5-N7-C8	-9.83	98.99	103.90
26	1H	729	G	C8-N9-C4	-9.82	102.47	106.40
26	1H	974(A)	C	N1-C2-O2	9.80	124.78	118.90
26	1H	793	A	N1-C6-N6	9.79	124.47	118.60
26	1H	621	A	C6-C5-N7	-9.76	125.47	132.30
26	1H	839	U	O5'-P-OP2	-9.76	96.92	105.70
26	14	2068	U	O5'-P-OP1	-9.76	96.92	105.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	1496	A	N7-C8-N9	9.76	118.68	113.80
26	1H	945	A	O4'-C1'-N9	9.74	116.00	108.20
26	14	2873	A	C2-N3-C4	-9.74	105.73	110.60
1	13	1158	C	N1-C2-O2	9.74	124.74	118.90
26	1H	1496	A	C4-C5-N7	9.74	115.57	110.70
26	1H	2085	C	O5'-P-OP2	-9.74	96.93	105.70
26	14	2502	G	O5'-P-OP1	-9.73	96.94	105.70
1	13	1502	A	C5-N7-C8	-9.71	99.05	103.90
26	14	2332	U	O5'-P-OP2	-9.70	96.97	105.70
26	1H	582	G	C5-C6-O6	-9.70	122.78	128.60
26	1H	1786	A	C4-C5-N7	9.68	115.54	110.70
26	1H	1816	G	O5'-P-OP1	-9.68	96.99	105.70
26	14	330	A	C2-N3-C4	-9.68	105.76	110.60
26	14	856	C	C6-N1-C2	-9.67	116.43	120.30
26	1H	837	C	N3-C4-N4	9.66	124.76	118.00
26	1H	1376	C	N3-C4-C5	-9.65	118.04	121.90
26	1H	2713	A	C2-N3-C4	-9.65	105.77	110.60
26	1H	793	A	C5-C6-N6	-9.65	115.98	123.70
26	14	565	C	C6-N1-C2	9.64	124.16	120.30
26	14	1681	G	C5-N7-C8	-9.64	99.48	104.30
26	14	2443	C	O5'-P-OP1	-9.64	97.02	105.70
26	1H	788	A	C5-C6-N1	-9.64	112.88	117.70
26	14	74	A	C5-C6-N1	-9.63	112.89	117.70
26	14	1332	G	C4-N9-C1'	9.63	139.01	126.50
26	1H	638	G	O5'-P-OP1	-9.62	97.04	105.70
26	14	1613	G	N3-C4-N9	9.62	131.77	126.00
26	1H	2346	A	C8-N9-C4	-9.62	101.95	105.80
26	1H	663	G	C8-N9-C4	-9.62	102.55	106.40
26	14	462	C	O5'-P-OP2	-9.61	97.05	105.70
26	1H	1772	G	C5-C6-O6	9.61	134.36	128.60
26	1H	667	U	N1-C2-O2	-9.60	116.08	122.80
26	14	140	A	N7-C8-N9	9.60	118.60	113.80
26	1H	245	G	C6-C5-N7	-9.60	124.64	130.40
26	1H	598	G	O5'-P-OP2	-9.59	97.07	105.70
26	1H	837	C	C5-C4-N4	-9.57	113.50	120.20
26	1H	1021	A	C2-N3-C4	-9.57	105.82	110.60
26	14	467	G	O5'-P-OP2	-9.56	97.09	105.70
26	1H	1255	U	N3-C4-O4	9.56	126.09	119.40
26	1H	2406	U	O5'-P-OP1	-9.56	97.10	105.70
26	14	1404	C	O5'-P-OP1	-9.55	97.10	105.70
26	14	2430	A	C6-C5-N7	-9.55	125.62	132.30
26	14	1902	C	O5'-P-OP1	-9.54	97.11	105.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	1006	C	O5'-P-OP1	-9.54	97.11	105.70
26	14	2873	A	C6-C5-N7	-9.54	125.63	132.30
26	14	2713	A	C5-N7-C8	-9.52	99.14	103.90
1	13	1502	A	O5'-P-OP2	-9.52	97.13	105.70
26	1H	676	A	N1-C2-N3	9.52	134.06	129.30
26	1H	783	A	N3-C4-C5	9.51	133.45	126.80
26	1H	2518	A	N7-C8-N9	9.50	118.55	113.80
1	1G	1517	G	O5'-P-OP2	-9.48	97.16	105.70
26	1H	1255	U	N3-C2-O2	9.48	128.84	122.20
26	14	2565	A	C8-N9-C4	9.47	109.59	105.80
26	1H	2287	A	C5-C6-N1	-9.47	112.96	117.70
26	14	49	A	O5'-P-OP2	-9.47	97.18	105.70
26	14	856	C	O5'-P-OP1	-9.47	97.18	105.70
26	1H	265	A	N7-C8-N9	9.46	118.53	113.80
26	1H	799	G	N7-C8-N9	-9.45	108.37	113.10
26	1H	1376	C	O5'-P-OP1	-9.45	97.19	105.70
1	13	690	G	O4'-C1'-N9	9.45	115.76	108.20
26	14	2873	A	N1-C2-N3	9.45	134.02	129.30
1	13	760	G	N1-C6-O6	9.44	125.57	119.90
26	1H	71	A	C4-C5-N7	9.44	115.42	110.70
26	14	2592	G	O5'-P-OP2	-9.44	97.21	105.70
26	1H	74	A	C5-C6-N1	-9.43	112.98	117.70
26	1H	2581	G	O5'-P-OP1	-9.43	97.21	105.70
26	14	621	A	C2-N3-C4	-9.43	105.88	110.60
26	1H	860	U	C2-N1-C1'	9.42	129.01	117.70
26	14	2276	G	O5'-P-OP1	-9.42	97.22	105.70
26	1H	491	G	O5'-P-OP1	-9.42	97.22	105.70
26	14	783	A	C6-C5-N7	-9.42	125.71	132.30
26	1H	1950	G	N7-C8-N9	9.41	117.80	113.10
26	14	2424	C	O5'-P-OP1	-9.39	97.25	105.70
1	13	1502	A	C4-C5-N7	9.38	115.39	110.70
26	1H	735	A	C8-N9-C4	9.37	109.55	105.80
26	1H	2392	A	N1-C6-N6	9.36	124.22	118.60
26	1H	757	U	O5'-P-OP2	-9.36	97.28	105.70
26	1H	71	A	N1-C2-N3	9.36	133.98	129.30
26	14	1678	G	N3-C4-C5	9.33	133.26	128.60
26	14	2056	G	C5-C6-O6	-9.33	123.00	128.60
26	1H	788	A	C6-N1-C2	9.32	124.19	118.60
26	1H	1779	U	OP1-P-OP2	9.32	133.59	119.60
27	16	30	C	C6-N1-C2	-9.32	116.57	120.30
26	1H	2249	U	O5'-P-OP1	-9.32	97.31	105.70
1	13	690	G	N1-C6-O6	9.31	125.48	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	1950	G	N3-C4-C5	-9.31	123.95	128.60
26	1H	733	G	OP1-P-OP2	9.30	133.56	119.60
24	3K	76	A	C8-N9-C4	-9.30	102.08	105.80
26	1H	930	U	C5-C4-O4	9.30	131.48	125.90
26	1H	593	G	O5'-P-OP2	-9.29	97.33	105.70
26	1H	1188	U	O5'-P-OP2	-9.29	97.34	105.70
26	1H	913	U	N3-C2-O2	-9.28	115.70	122.20
26	1H	2346	A	N7-C8-N9	9.28	118.44	113.80
26	1H	774	A	C4-C5-C6	-9.28	112.36	117.00
26	14	1673	U	O5'-P-OP1	-9.28	97.35	105.70
26	1H	2705	A	N1-C6-N6	9.27	124.16	118.60
1	13	281	G	O5'-P-OP1	-9.26	97.36	105.70
26	1H	512	G	O4'-C1'-N9	9.26	115.60	108.20
26	14	2430	A	C4-C5-C6	9.26	121.63	117.00
26	1H	774	A	C8-N9-C4	-9.25	102.10	105.80
26	1H	1678	G	C8-N9-C4	-9.25	102.70	106.40
26	1H	2330	G	C5-C6-O6	-9.23	123.06	128.60
26	1H	2392	A	C2-N3-C4	-9.23	105.99	110.60
26	1H	2374	C	C6-N1-C2	9.22	123.99	120.30
26	14	801	G	N1-C6-O6	-9.22	114.37	119.90
26	14	1616	A	C2-N3-C4	-9.20	106.00	110.60
1	13	811	C	C6-N1-C2	9.20	123.98	120.30
26	1H	2507	C	C6-N1-C2	-9.20	116.62	120.30
26	1H	728	G	O5'-P-OP2	-9.18	97.44	105.70
26	1H	203	C	N1-C2-O2	-9.18	113.39	118.90
26	14	2870	C	C6-N1-C2	-9.18	116.63	120.30
26	1H	1899	G	C8-N9-C4	-9.18	102.73	106.40
26	1H	2747	G	C6-C5-N7	-9.17	124.90	130.40
1	13	1266	G	N3-C4-N9	-9.17	120.50	126.00
26	1H	1558	A	O5'-P-OP1	-9.17	97.45	105.70
26	14	530	G	C6-C5-N7	-9.17	124.90	130.40
26	14	2253	G	N1-C6-O6	9.17	125.40	119.90
26	1H	2688	U	C5-C4-O4	9.16	131.40	125.90
26	14	2438	U	O5'-P-OP2	-9.16	97.45	105.70
26	14	2518	A	C6-C5-N7	-9.16	125.89	132.30
26	1H	2582	G	N3-C4-N9	9.16	131.49	126.00
26	14	252	G	N1-C6-O6	-9.13	114.42	119.90
26	14	1359	A	C8-N9-C4	9.13	109.45	105.80
26	14	209	C	C5-C4-N4	-9.12	113.81	120.20
26	1H	621	A	N1-C2-N3	9.11	133.86	129.30
26	14	774	A	O5'-P-OP2	-9.11	97.50	105.70
1	13	789	U	N3-C2-O2	-9.11	115.83	122.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	2212	A	O4'-C1'-N9	9.10	115.48	108.20
26	1H	2585	U	N1-C2-O2	9.10	129.17	122.80
26	1H	1574	C	C6-N1-C2	9.10	123.94	120.30
36	35	147	LEU	CA-CB-CG	9.10	136.22	115.30
1	13	529	G	N1-C6-O6	9.09	125.36	119.90
26	1H	2010	G	O5'-P-OP1	-9.09	97.52	105.70
26	1H	1255	U	N1-C2-O2	-9.08	116.44	122.80
26	1H	2070	G	N1-C6-O6	-9.08	114.45	119.90
39	65	110	LEU	CA-CB-CG	9.08	136.19	115.30
26	1H	1300	U	N1-C2-N3	9.08	120.35	114.90
26	14	729	G	N3-C2-N2	-9.08	113.54	119.90
26	1H	1306	C	C6-N1-C2	9.08	123.93	120.30
26	14	2001	A	N1-C6-N6	9.08	124.05	118.60
26	1H	528	A	N3-C4-C5	9.07	133.15	126.80
26	1H	1990	C	C6-N1-C2	-9.07	116.67	120.30
26	14	1340	U	C6-N1-C2	9.07	126.44	121.00
27	16	79	C	C6-N1-C2	-9.07	116.67	120.30
26	1H	1332	G	N3-C4-N9	-9.06	120.57	126.00
26	1H	578	A	O5'-P-OP2	-9.06	97.55	105.70
34	15	42	TRP	CE2-CD2-CE3	9.05	129.57	118.70
26	1H	2380	C	C2-N3-C4	-9.05	115.37	119.90
26	14	140	A	C4-C5-N7	9.04	115.22	110.70
26	14	1325	G	C5-C6-O6	-9.04	123.17	128.60
26	1H	1386	C	C6-N1-C2	-9.04	116.68	120.30
26	14	2070	G	O5'-P-OP2	-9.02	97.58	105.70
26	1H	2503	A	N1-C2-N3	-9.01	124.79	129.30
26	1H	2584	U	C5-C4-O4	9.01	131.31	125.90
26	1H	1617	C	O5'-P-OP1	-9.00	97.60	105.70
26	1H	71	A	N7-C8-N9	8.99	118.30	113.80
1	13	1502	A	N1-C6-N6	8.99	123.99	118.60
1	13	792	A	N1-C6-N6	8.97	123.98	118.60
26	1H	138	G	C8-N9-C4	-8.97	102.81	106.40
26	14	250	G	O5'-P-OP1	-8.97	97.63	105.70
26	1H	71	A	O4'-C1'-N9	-8.97	101.03	108.20
26	1H	567	A	O5'-P-OP1	-8.96	97.63	105.70
26	1H	641	C	O5'-P-OP1	-8.96	97.64	105.70
26	1H	1382	G	N1-C6-O6	8.96	125.28	119.90
26	1H	1193	G	O5'-P-OP2	-8.96	97.64	105.70
26	1H	1814	G	N3-C4-N9	8.95	131.37	126.00
26	14	2056	G	N3-C2-N2	-8.94	113.64	119.90
26	1H	691	C	C6-N1-C2	8.94	123.88	120.30
26	14	1661	G	C8-N9-C4	8.94	109.98	106.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	2468	G	O4'-C1'-N9	8.93	115.35	108.20
26	1H	1814	G	C5-C6-N1	8.93	115.96	111.50
26	1H	245	G	C5-C6-O6	-8.93	123.25	128.60
26	14	774	A	C4-C5-C6	-8.92	112.54	117.00
26	1H	1528	A	N7-C8-N9	8.92	118.26	113.80
26	14	791	C	C6-N1-C2	8.92	123.87	120.30
26	1H	2688	U	N3-C2-O2	-8.90	115.97	122.20
26	14	774	A	C2-N3-C4	-8.90	106.15	110.60
26	1H	1698	A	C5-N7-C8	-8.89	99.45	103.90
26	14	2518	A	C4-C5-N7	8.89	115.14	110.70
1	13	792	A	C6-C5-N7	-8.89	126.08	132.30
26	1H	530	G	C5-C6-O6	8.88	133.93	128.60
26	1H	2298	A	O5'-P-OP2	-8.88	97.71	105.70
55	Q8	25	MET	N-CA-C	8.88	134.98	111.00
1	1G	14	U	C5-C6-N1	8.88	127.14	122.70
26	14	1902	C	N3-C4-C5	8.88	125.45	121.90
26	1H	1814	G	N9-C4-C5	-8.87	101.85	105.40
26	14	2542	A	C8-N9-C4	8.86	109.34	105.80
26	1H	140	A	C6-C5-N7	-8.85	126.10	132.30
1	13	1260	C	C6-N1-C2	-8.85	116.76	120.30
26	1H	2712	U	C5-C6-N1	-8.84	118.28	122.70
1	13	745	C	C6-N1-C2	-8.84	116.77	120.30
26	14	821	A	N1-C6-N6	8.83	123.90	118.60
26	14	1786	A	C4-C5-N7	8.82	115.11	110.70
26	14	769	G	OP1-P-O3'	8.81	124.59	105.20
26	14	1496	A	C8-N9-C4	-8.80	102.28	105.80
26	1H	2311	A	N1-C2-N3	8.79	133.69	129.30
26	1H	2228	G	C6-C5-N7	-8.79	125.13	130.40
26	14	2272	U	N3-C2-O2	-8.79	116.05	122.20
23	2K	1	C	C6-N1-C2	-8.78	116.79	120.30
26	1H	1828	G	C5-C6-N1	-8.78	107.11	111.50
22	1K	76	A	C8-N9-C4	-8.77	102.29	105.80
26	1H	2270	G	C8-N9-C4	8.77	109.91	106.40
26	14	71	A	C2-N3-C4	-8.76	106.22	110.60
26	1H	1577	C	N3-C4-C5	-8.75	118.40	121.90
26	14	2501	C	C2-N1-C1'	-8.74	109.18	118.80
26	14	1774	C	O5'-P-OP1	-8.74	97.84	105.70
26	14	2512	C	N3-C4-C5	8.74	125.39	121.90
26	1H	1204	A	C2-N3-C4	-8.73	106.23	110.60
26	14	737	C	N1-C2-O2	-8.73	113.66	118.90
26	1H	693	C	C5-C4-N4	8.73	126.31	120.20
26	1H	1379	A	C5-N7-C8	-8.73	99.53	103.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	528	A	N3-C4-N9	-8.73	120.42	127.40
26	1H	189	G	C8-N9-C4	8.73	109.89	106.40
26	1H	140	A	C4-C5-N7	8.72	115.06	110.70
26	1H	144	C	C6-N1-C2	8.72	123.79	120.30
26	1H	463	G	O5'-P-OP2	-8.72	97.85	105.70
26	14	1827	C	OP1-P-O3'	8.72	124.39	105.20
1	13	974	A	N1-C6-N6	8.72	123.83	118.60
26	1H	691	C	N1-C2-O2	-8.71	113.67	118.90
26	1H	739	G	C8-N9-C4	8.71	109.88	106.40
26	1H	1616	A	C4-C5-N7	8.71	115.05	110.70
26	14	676	A	O4'-C1'-N9	8.69	115.15	108.20
26	1H	690	G	C8-N9-C4	8.69	109.88	106.40
26	1H	1517	G	OP1-P-O3'	8.69	124.31	105.20
26	14	1698	A	C4-C5-N7	8.69	115.04	110.70
26	14	1786	A	C6-C5-N7	-8.68	126.22	132.30
26	1H	858	U	O5'-P-OP2	-8.66	97.90	105.70
26	1H	2508	G	N9-C4-C5	8.66	108.87	105.40
26	1H	2053	G	N1-C6-O6	8.65	125.09	119.90
26	14	2307	G	O4'-C1'-N9	8.65	115.12	108.20
1	13	542	G	O5'-P-OP1	-8.65	97.92	105.70
26	14	1279	G	N1-C6-O6	-8.65	114.71	119.90
26	1H	1786	A	C5-C6-N1	-8.64	113.38	117.70
26	1H	2598	A	O5'-P-OP2	8.64	121.07	110.70
26	1H	957	A	N1-C6-N6	8.63	123.78	118.60
26	1H	1828	G	O5'-P-OP2	-8.64	97.93	105.70
26	1H	2331	G	C8-N9-C4	8.63	109.85	106.40
26	1H	2236	C	O5'-P-OP1	-8.62	97.94	105.70
26	1H	398	G	O5'-P-OP2	-8.62	97.94	105.70
26	1H	211	A	C8-N9-C4	8.62	109.25	105.80
26	1H	964	C	O5'-P-OP1	-8.62	97.94	105.70
26	1H	138	G	N7-C8-N9	8.62	117.41	113.10
26	14	2287	A	C2-N3-C4	-8.61	106.29	110.60
26	1H	74	A	C5-N7-C8	-8.61	99.60	103.90
26	14	1678	G	N3-C4-N9	-8.61	120.83	126.00
26	14	1908	C	C6-N1-C2	-8.61	116.86	120.30
26	1H	746	A	O5'-P-OP2	8.60	121.03	110.70
26	1H	226	G	O4'-C1'-N9	8.60	115.08	108.20
26	1H	752	A	P-O3'-C3'	8.60	130.02	119.70
26	1H	1610	A	C2-N3-C4	-8.60	106.30	110.60
26	1H	2330	G	N1-C6-O6	8.60	125.06	119.90
26	1H	1379	A	C4-C5-N7	8.60	115.00	110.70
26	1H	245	G	N1-C6-O6	8.59	125.05	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	2054	A	OP2-P-O3'	8.58	124.08	105.20
26	1H	391	G	C6-C5-N7	-8.57	125.26	130.40
26	1H	774	A	C8-N9-C1'	8.57	143.13	127.70
26	1H	613	U	C5-C4-O4	8.57	131.04	125.90
26	14	829	A	OP1-P-OP2	8.57	132.45	119.60
26	1H	774	A	N1-C2-N3	-8.56	125.02	129.30
26	1H	1310	G	O5'-P-OP2	8.56	120.97	110.70
26	1H	617	G	C8-N9-C4	8.56	109.82	106.40
26	14	1342	A	N1-C6-N6	8.56	123.74	118.60
26	1H	1600	C	O5'-P-OP2	-8.56	98.00	105.70
26	14	1907	G	O5'-P-OP1	-8.56	98.00	105.70
26	1H	1820	U	C5-C6-N1	-8.55	118.42	122.70
37	88	82	ARG	N-CA-C	8.55	134.08	111.00
26	1H	2499	C	N1-C2-O2	-8.54	113.78	118.90
26	1H	245	G	C4-C5-N7	8.53	114.21	110.80
26	14	140	A	N1-C6-N6	8.53	123.72	118.60
26	1H	141	A	C5-N7-C8	-8.53	99.64	103.90
26	1H	1899	G	C5-C6-N1	-8.53	107.24	111.50
26	14	1786	A	C8-N9-C4	-8.52	102.39	105.80
26	14	2426	A	N7-C8-N9	8.52	118.06	113.80
26	1H	679	C	C6-N1-C2	8.51	123.70	120.30
26	1H	429	A	C8-N9-C4	-8.51	102.40	105.80
26	1H	796	C	C6-N1-C2	8.51	123.70	120.30
26	14	929	G	C6-C5-N7	-8.51	125.30	130.40
26	1H	2330	G	C6-C5-N7	-8.50	125.30	130.40
26	14	1394	U	O5'-P-OP2	8.50	120.90	110.70
26	1H	1431	U	C5-C6-N1	8.50	126.95	122.70
26	1H	2518	A	C4-C5-N7	8.49	114.95	110.70
26	1H	2451	A	N1-C6-N6	-8.49	113.51	118.60
23	2L	21	U	N1-C2-N3	8.49	119.99	114.90
26	14	2873	A	C8-N9-C4	-8.49	102.40	105.80
26	1H	1428	C	C6-N1-C2	8.49	123.69	120.30
26	1H	1528	A	O4'-C1'-N9	8.49	114.99	108.20
26	14	1616	A	N7-C8-N9	8.49	118.04	113.80
26	1H	1616	A	O4'-C1'-N9	8.48	114.98	108.20
26	1H	2425	A	O5'-P-OP2	-8.47	98.07	105.70
26	14	1618	A	C8-N9-C4	-8.47	102.41	105.80
26	1H	828	U	N1-C2-O2	8.46	128.72	122.80
26	1H	1660	C	N3-C2-O2	-8.46	115.97	121.90
23	2K	17	C	N1-C2-O2	8.46	123.98	118.90
26	1H	1636	C	N3-C4-C5	-8.45	118.52	121.90
26	14	463	G	O5'-P-OP2	-8.45	98.10	105.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	13	812	C	N1-C2-O2	8.44	123.97	118.90
36	78	20	GLY	N-CA-C	8.44	134.20	113.10
26	14	1762	A	O4'-C1'-N9	8.43	114.94	108.20
26	14	2873	A	N1-C6-N6	8.43	123.66	118.60
26	1H	1784	A	O4'-C1'-N9	-8.42	101.46	108.20
26	1H	2010	G	O5'-P-OP2	8.42	120.81	110.70
26	1H	2747	G	C5-C6-O6	-8.42	123.55	128.60
1	1G	254	G	O5'-P-OP1	-8.42	98.13	105.70
26	14	1898	U	N3-C4-C5	-8.42	109.55	114.60
26	14	1407	C	C5-C6-N1	8.41	125.21	121.00
26	14	1342	A	C2-N3-C4	-8.41	106.39	110.60
26	1H	330	A	N1-C2-N3	8.41	133.50	129.30
26	1H	788	A	N9-C4-C5	-8.40	102.44	105.80
26	14	1396	U	N3-C2-O2	-8.40	116.32	122.20
26	1H	2582	G	N9-C4-C5	-8.39	102.04	105.40
26	1H	1598	C	OP1-P-O3'	8.39	123.65	105.20
26	1H	1210	A	N1-C6-N6	8.38	123.63	118.60
26	14	1340	U	C5-C6-N1	-8.38	118.51	122.70
26	14	2087	G	O5'-P-OP2	-8.38	98.16	105.70
26	1H	972	G	N1-C6-O6	-8.37	114.88	119.90
26	1H	2392	A	N7-C8-N9	8.37	117.98	113.80
26	14	774	A	C6-N1-C2	8.37	123.62	118.60
26	1H	528	A	C5-C6-N1	-8.36	113.52	117.70
26	1H	2699	C	N3-C4-C5	8.36	125.25	121.90
26	1H	2430	A	C5-C6-N1	-8.36	113.52	117.70
26	1H	659	C	C6-N1-C2	8.36	123.64	120.30
26	14	676	A	N3-C4-C5	8.35	132.65	126.80
26	1H	1779	U	O5'-P-OP1	-8.35	98.19	105.70
26	14	2078	C	N3-C4-C5	-8.35	118.56	121.90
26	1H	2377	A	C8-N9-C4	8.35	109.14	105.80
26	14	737	C	N3-C4-N4	8.34	123.84	118.00
26	1H	2272	U	O5'-P-OP1	8.34	120.70	110.70
26	14	2429	G	O5'-P-OP2	-8.34	98.20	105.70
26	1H	2433	A	N1-C2-N3	8.33	133.47	129.30
26	14	2324	C	C6-N1-C2	8.33	123.63	120.30
26	1H	783	A	C5-C6-N1	-8.33	113.54	117.70
26	14	729	G	N1-C2-N2	8.33	123.69	116.20
23	2L	21	U	C5-C4-O4	-8.32	120.91	125.90
26	14	1698	A	C5-N7-C8	-8.32	99.74	103.90
30	31	176	LEU	CA-CB-CG	8.32	134.44	115.30
26	1H	634	C	O5'-P-OP2	-8.32	98.21	105.70
26	1H	961	C	C6-N1-C2	-8.31	116.97	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	1142(A)	A	C2-N3-C4	-8.31	106.44	110.60
26	14	1327	C	N1-C2-O2	-8.31	113.91	118.90
26	1H	471	A	C2-N3-C4	-8.31	106.44	110.60
26	1H	140	A	N1-C6-N6	8.31	123.58	118.60
26	1H	681	G	N1-C2-N3	8.30	128.88	123.90
26	14	1678	G	N7-C8-N9	8.30	117.25	113.10
26	14	1394	U	O5'-P-OP1	-8.30	98.23	105.70
26	1H	1021	A	C5-N7-C8	-8.29	99.75	103.90
26	1H	459	U	O5'-P-OP2	-8.29	98.24	105.70
26	14	74	A	N3-C4-C5	8.29	132.60	126.80
26	14	2518	A	C2-N3-C4	-8.29	106.46	110.60
26	1H	2782	G	N1-C6-O6	8.28	124.87	119.90
26	1H	25	U	C5-C4-O4	-8.28	120.94	125.90
26	1H	2072	G	OP1-P-O3'	8.28	123.41	105.20
26	1H	2575	C	C5-C6-N1	-8.28	116.86	121.00
26	1H	2392	A	C4-C5-N7	8.27	114.84	110.70
26	14	818	G	C8-N9-C4	-8.27	103.09	106.40
26	1H	794	G	O5'-P-OP1	-8.27	98.26	105.70
26	14	1613	G	N3-C2-N2	8.27	125.69	119.90
26	1H	144	C	N3-C4-C5	8.27	125.21	121.90
55	Q8	60	LEU	CA-CB-CG	-8.27	96.29	115.30
26	1H	2822	G	N1-C6-O6	8.26	124.86	119.90
26	14	1681	G	C4-C5-N7	8.26	114.11	110.80
24	3K	76	A	C4-C5-N7	8.26	114.83	110.70
26	1H	2490	G	O5'-P-OP2	-8.26	98.27	105.70
1	13	811	C	C5-C6-N1	-8.26	116.87	121.00
1	13	1502	A	C2-N3-C4	-8.25	106.47	110.60
24	3L	76	A	C8-N9-C4	-8.25	102.50	105.80
1	13	1482	G	O5'-P-OP2	-8.24	98.28	105.70
26	1H	774	A	C2-N3-C4	-8.24	106.48	110.60
26	14	558	G	C8-N9-C4	8.24	109.70	106.40
26	1H	913	U	O5'-P-OP2	-8.23	98.29	105.70
26	1H	1567	A	N1-C6-N6	-8.23	113.66	118.60
26	1H	1629	U	O5'-P-OP2	8.23	120.57	110.70
34	15	42	TRP	CE2-CD2-CG	-8.23	100.72	107.30
26	1H	821	A	OP1-P-OP2	8.22	131.94	119.60
34	15	42	TRP	CD2-CE2-CZ2	-8.22	112.43	122.30
26	1H	769	G	N1-C2-N2	-8.21	108.81	116.20
1	1G	117	G	N1-C6-O6	8.21	124.83	119.90
26	1H	121	G	C5-C6-O6	-8.21	123.68	128.60
26	1H	141	A	N7-C8-N9	8.20	117.90	113.80
26	14	1253	A	C5-C6-N6	-8.21	117.14	123.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	2880	C	C6-N1-C2	-8.20	117.02	120.30
26	1H	1950	G	C4-C5-N7	8.20	114.08	110.80
26	14	1950	G	C4-N9-C1'	8.20	137.16	126.50
1	13	760	G	C6-C5-N7	-8.20	125.48	130.40
26	1H	144	C	C5-C6-N1	-8.19	116.90	121.00
26	14	1614	A	C2-N3-C4	-8.19	106.50	110.60
26	1H	945	A	C5-C6-N1	-8.19	113.61	117.70
26	1H	245	G	N9-C4-C5	-8.19	102.12	105.40
26	14	1613	G	N9-C4-C5	-8.19	102.12	105.40
26	14	2357	U	O5'-P-OP2	-8.19	98.33	105.70
1	13	1496	C	C2-N1-C1'	-8.18	109.80	118.80
26	1H	705	A	N1-C6-N6	8.18	123.51	118.60
26	1H	2346	A	C5-N7-C8	-8.18	99.81	103.90
26	1H	1785	A	C4-C5-C6	8.18	121.09	117.00
26	14	1496	A	C5-N7-C8	-8.18	99.81	103.90
26	1H	1626	G	C8-N9-C4	-8.17	103.13	106.40
26	14	2713	A	N7-C8-N9	8.17	117.89	113.80
26	14	2253	G	C5-C6-O6	-8.16	123.70	128.60
26	14	2595	G	O5'-P-OP1	-8.16	98.36	105.70
26	14	1145	C	C6-N1-C2	-8.15	117.04	120.30
26	1H	123	G	O5'-P-OP2	-8.15	98.37	105.70
27	16	79	C	C2-N1-C1'	8.15	127.76	118.80
1	13	1502	A	C6-C5-N7	-8.14	126.60	132.30
26	1H	2702	U	C5-C6-N1	8.14	126.77	122.70
26	14	1520	U	C5-C4-O4	8.14	130.78	125.90
26	14	2437	U	C5-C4-O4	8.14	130.78	125.90
26	1H	263	C	O5'-P-OP2	-8.13	98.38	105.70
26	14	921	G	C8-N9-C4	-8.13	103.15	106.40
26	1H	2430	A	N3-C4-N9	-8.13	120.90	127.40
26	1H	189	G	O5'-P-OP2	8.13	120.45	110.70
26	1H	1350	C	C6-N1-C2	8.12	123.55	120.30
26	1H	1314	C	C2-N1-C1'	8.12	127.73	118.80
26	1H	2259	G	OP1-P-OP2	-8.12	107.43	119.60
26	14	2688	U	C5-C6-N1	-8.11	118.64	122.70
26	1H	1526	G	C8-N9-C4	-8.11	103.16	106.40
26	1H	801	G	N3-C2-N2	-8.11	114.23	119.90
26	14	2335	A	O4'-C1'-N9	8.10	114.68	108.20
26	1H	1496	A	N1-C6-N6	8.10	123.46	118.60
26	1H	146	G	C4-C5-N7	8.09	114.04	110.80
1	13	690	G	C4-C5-C6	8.09	123.66	118.80
26	1H	2446	G	C4-C5-N7	8.09	114.04	110.80
26	14	2518	A	O4'-C1'-N9	-8.09	101.73	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	13	808	C	N3-C4-C5	-8.09	118.67	121.90
26	1H	820	A	O5'-P-OP1	-8.09	98.42	105.70
26	14	676	A	N7-C8-N9	8.08	117.84	113.80
26	1H	189	G	N1-C6-O6	8.08	124.75	119.90
26	1H	1201	C	N3-C2-O2	8.08	127.56	121.90
26	1H	265	A	C5-N7-C8	-8.08	99.86	103.90
26	1H	2346	A	C6-C5-N7	-8.07	126.65	132.30
27	16	101	A	O5'-P-OP2	-8.07	98.43	105.70
1	1G	449	C	C6-N1-C2	-8.07	117.07	120.30
26	14	1970	A	O5'-P-OP2	-8.07	98.44	105.70
26	1H	681	G	N1-C2-N2	-8.07	108.94	116.20
26	14	71	A	C5-N7-C8	-8.07	99.86	103.90
26	1H	130	C	C5-C4-N4	-8.07	114.55	120.20
26	1H	1401	G	C8-N9-C4	-8.07	103.17	106.40
26	14	570	G	N3-C4-C5	-8.07	124.57	128.60
26	1H	462	C	C6-N1-C2	-8.06	117.08	120.30
26	14	783	A	C5-C6-N1	-8.06	113.67	117.70
26	1H	321	G	C6-C5-N7	-8.06	125.57	130.40
23	2K	13	C	C5-C6-N1	8.06	125.03	121.00
26	1H	659	C	C5-C6-N1	-8.06	116.97	121.00
26	1H	2688	U	N1-C2-N3	8.05	119.73	114.90
26	1H	2747	G	C4-C5-N7	8.05	114.02	110.80
26	1H	2577	A	N1-C6-N6	-8.05	113.77	118.60
26	1H	1518	C	O5'-P-OP1	-8.05	98.46	105.70
26	1H	2582	G	C4-C5-N7	8.05	114.02	110.80
26	1H	929	G	N1-C6-O6	8.04	124.73	119.90
26	1H	2507	C	N1-C2-O2	8.05	123.73	118.90
26	14	2249	U	C6-N1-C2	-8.04	116.18	121.00
26	1H	774	A	N7-C8-N9	8.04	117.82	113.80
26	14	121	G	C5-C6-O6	-8.04	123.78	128.60
26	1H	2282	G	O5'-P-OP1	-8.04	98.47	105.70
26	14	2351	G	N3-C4-C5	-8.04	124.58	128.60
26	1H	2518	A	C8-N9-C4	-8.03	102.59	105.80
26	1H	2252	G	C8-N9-C4	8.03	109.61	106.40
28	19	272	ALA	N-CA-C	8.03	132.67	111.00
26	1H	2238	G	O5'-P-OP2	-8.02	98.48	105.70
26	1H	2232	U	C5-C4-O4	8.02	130.71	125.90
26	1H	1950	G	O4'-C1'-N9	8.02	114.61	108.20
26	1H	1254	A	C2-N3-C4	-8.02	106.59	110.60
26	1H	2819	G	C5-C6-O6	-8.01	123.79	128.60
1	13	362	G	O5'-P-OP1	-8.01	98.50	105.70
1	13	528	C	C6-N1-C2	8.01	123.50	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	1612	C	C6-N1-C2	8.01	123.50	120.30
26	1H	1022	G	N9-C4-C5	8.00	108.60	105.40
26	1H	1192	G	C8-N9-C4	8.00	109.60	106.40
22	1K	76	A	N7-C8-N9	7.99	117.80	113.80
26	1H	1781	C	N3-C2-O2	-7.99	116.31	121.90
26	1H	2435	A	O5'-P-OP1	-7.99	98.51	105.70
26	1H	784	A	O4'-C1'-N9	7.98	114.59	108.20
26	1H	2439	A	O5'-P-OP2	-7.98	98.52	105.70
26	14	2443	C	O5'-P-OP2	7.98	120.28	110.70
26	1H	1639	U	N3-C2-O2	-7.98	116.62	122.20
26	14	1324	G	N1-C6-O6	7.98	124.69	119.90
1	13	254	G	O5'-P-OP1	-7.98	98.52	105.70
26	1H	1605	C	O5'-P-OP1	-7.98	98.52	105.70
26	14	1342	A	N1-C2-N3	7.98	133.29	129.30
26	1H	286	C	O5'-P-OP2	-7.97	98.52	105.70
26	1H	222	A	P-O3'-C3'	7.97	129.27	119.70
26	14	2610	C	C5-C4-N4	-7.97	114.62	120.20
26	1H	917	A	N1-C6-N6	7.96	123.38	118.60
26	1H	797	C	C6-N1-C2	7.96	123.48	120.30
26	1H	2392	A	C5-C6-N1	-7.96	113.72	117.70
26	14	1932	A	O5'-P-OP1	-7.96	98.54	105.70
26	1H	1931	U	C5-C6-N1	-7.95	118.72	122.70
26	1H	2267	A	OP1-P-O3'	7.95	122.69	105.20
26	1H	2698	U	C5-C6-N1	-7.95	118.73	122.70
26	1H	1324	G	N1-C6-O6	7.94	124.67	119.90
26	1H	1520	U	N3-C2-O2	-7.94	116.64	122.20
26	14	929	G	N7-C8-N9	7.94	117.07	113.10
26	14	1306	C	O5'-P-OP1	-7.93	98.56	105.70
26	1H	2248	C	N3-C4-N4	-7.92	112.45	118.00
26	14	774	A	O5'-P-OP1	7.92	120.21	110.70
26	14	792	G	N3-C4-C5	-7.92	124.64	128.60
26	1H	834	C	OP2-P-O3'	7.92	122.63	105.20
26	1H	865	C	O5'-P-OP2	7.92	120.21	110.70
26	1H	1314	C	O5'-P-OP2	-7.92	98.57	105.70
26	1H	1496	A	C8-N9-C4	-7.92	102.63	105.80
1	13	353	A	C8-N9-C4	-7.92	102.63	105.80
26	1H	48	G	OP2-P-O3'	7.91	122.61	105.20
26	1H	692	C	C5-C4-N4	-7.91	114.66	120.20
26	1H	860	U	N1-C2-O2	7.91	128.34	122.80
1	13	789	U	C4-C5-C6	7.91	124.45	119.70
26	14	967	C	O5'-P-OP2	-7.91	98.58	105.70
26	1H	2576	G	C8-N9-C4	7.91	109.56	106.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	429	A	O5'-P-OP1	-7.90	98.59	105.70
26	14	1617	C	N1-C2-O2	-7.90	114.16	118.90
26	1H	2572	A	O5'-P-OP2	-7.90	98.59	105.70
26	14	676	A	C4-C5-N7	7.90	114.65	110.70
26	14	1283	G	N3-C4-C5	-7.90	124.65	128.60
26	1H	1021	A	N7-C8-N9	7.89	117.75	113.80
26	14	2439	A	N7-C8-N9	7.88	117.74	113.80
26	1H	946	G	C8-N9-C4	7.88	109.55	106.40
26	14	2542	A	N7-C8-N9	-7.87	109.86	113.80
26	1H	1302	A	OP1-P-OP2	7.87	131.40	119.60
26	1H	510	C	O5'-P-OP2	-7.86	98.62	105.70
26	1H	908	C	O5'-P-OP2	-7.86	98.62	105.70
26	14	945	A	C5-C6-N6	-7.86	117.41	123.70
1	13	990	C	C6-N1-C2	-7.86	117.16	120.30
26	1H	1622	G	N3-C2-N2	-7.86	114.40	119.90
26	1H	683	C	N3-C4-C5	7.85	125.04	121.90
36	35	62	LEU	N-CA-C	7.85	132.20	111.00
26	1H	693	C	N3-C4-N4	-7.85	112.51	118.00
26	14	1253	A	N9-C4-C5	-7.85	102.66	105.80
26	1H	49	A	C5-N7-C8	7.84	107.82	103.90
26	1H	942	G	N3-C2-N2	-7.84	114.41	119.90
26	1H	777	A	N1-C2-N3	7.84	133.22	129.30
26	1H	2457	U	N3-C2-O2	7.84	127.69	122.20
26	1H	1157	G	N1-C6-O6	7.83	124.60	119.90
26	1H	2518	A	N1-C6-N6	7.83	123.30	118.60
28	11	238	GLY	N-CA-C	7.83	132.67	113.10
1	1G	1397	C	C6-N1-C2	-7.83	117.17	120.30
26	1H	2380	C	C6-N1-C2	7.82	123.43	120.30
26	1H	664	C	O5'-P-OP2	-7.82	98.66	105.70
26	14	2046	G	N1-C6-O6	-7.82	115.21	119.90
26	14	2444	G	N1-C6-O6	-7.82	115.21	119.90
1	13	1322	C	C5-C6-N1	7.81	124.91	121.00
26	1H	839	U	OP1-P-OP2	7.81	131.31	119.60
42	D8	40	LEU	CA-CB-CG	7.81	133.26	115.30
26	1H	1526	G	N7-C8-N9	7.80	117.00	113.10
26	1H	640	C	O5'-P-OP2	-7.80	98.68	105.70
26	1H	1443	G	C4-C5-C6	7.80	123.48	118.80
26	14	783	A	N3-C4-N9	-7.80	121.16	127.40
27	16	79	C	C5-C6-N1	7.79	124.89	121.00
26	14	2267	A	OP1-P-OP2	7.79	131.29	119.60
26	1H	1210	A	C4-C5-N7	7.79	114.59	110.70
26	1H	529	A	N7-C8-N9	7.78	117.69	113.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	1G	1346	A	P-O3'-C3'	7.78	129.04	119.70
26	14	2617	C	C6-N1-C2	7.78	123.41	120.30
1	13	792	A	C2-N3-C4	-7.78	106.71	110.60
25	4K	18	G	N9-C4-C5	7.78	108.51	105.40
26	1H	526	A	N1-C6-N6	-7.78	113.93	118.60
26	1H	2003	G	O5'-P-OP1	-7.78	98.70	105.70
26	14	741	G	O5'-P-OP2	7.78	120.03	110.70
1	13	898	G	O5'-P-OP1	-7.78	98.70	105.70
26	1H	758	C	N3-C4-C5	7.78	125.01	121.90
26	1H	783	A	N1-C2-N3	7.77	133.18	129.30
26	1H	1888	G	N3-C4-C5	-7.76	124.72	128.60
26	1H	768	G	O5'-P-OP2	-7.76	98.72	105.70
26	14	1313	U	C2-N1-C1'	7.76	127.02	117.70
26	1H	863	A	O5'-P-OP1	7.76	120.01	110.70
23	2L	21	U	N3-C4-C5	7.76	119.26	114.60
26	14	2731	G	C6-C5-N7	-7.76	125.75	130.40
1	13	1468	A	C8-N9-C4	7.76	108.90	105.80
26	1H	636	G	O5'-P-OP2	7.75	120.00	110.70
26	14	2051	A	C8-N9-C4	-7.75	102.70	105.80
26	1H	119	A	C5-N7-C8	7.75	107.77	103.90
26	14	746	A	O5'-P-OP2	7.75	119.99	110.70
1	13	1158	C	C2-N1-C1'	7.74	127.32	118.80
26	14	621	A	C5-C6-N1	-7.74	113.83	117.70
26	14	1925	C	N1-C2-O2	-7.74	114.25	118.90
26	14	2518	A	N7-C8-N9	7.74	117.67	113.80
26	1H	1225	C	C6-N1-C2	7.74	123.40	120.30
26	1H	146	G	C5-C6-O6	-7.74	123.96	128.60
26	1H	528	A	C2-N3-C4	-7.72	106.74	110.60
26	14	1332	G	C8-N9-C1'	-7.72	116.96	127.00
1	1G	337	C	C6-N1-C2	-7.72	117.21	120.30
26	1H	537	C	O5'-P-OP1	7.72	119.96	110.70
26	14	561	G	N3-C4-N9	-7.72	121.37	126.00
23	2L	21	U	C2-N3-C4	-7.71	122.37	127.00
26	1H	739	G	N7-C8-N9	-7.71	109.24	113.10
26	1H	945	A	C8-N9-C1'	-7.71	113.82	127.70
26	1H	2228	G	C4-N9-C1'	7.71	136.53	126.50
26	1H	1300	U	N1-C2-O2	-7.71	117.41	122.80
1	13	827	U	N3-C2-O2	-7.70	116.81	122.20
26	1H	1428	C	N3-C4-N4	-7.70	112.61	118.00
26	1H	681	G	C2-N3-C4	-7.69	108.06	111.90
26	1H	1336	A	N9-C4-C5	7.69	108.87	105.80
26	14	1671	U	O5'-P-OP1	-7.69	98.78	105.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	1G	811	C	N1-C2-O2	-7.68	114.29	118.90
26	14	1527	G	N3-C4-N9	-7.68	121.39	126.00
26	1H	1614	A	N3-C4-C5	7.68	132.18	126.80
29	29	23	VAL	N-CA-C	7.68	131.74	111.00
26	1H	1611	C	C5-C6-N1	-7.67	117.16	121.00
26	1H	1022	G	C4-C5-N7	-7.67	107.73	110.80
1	13	816	A	C8-N9-C4	-7.67	102.73	105.80
26	1H	1297	C	O5'-P-OP2	-7.67	98.80	105.70
26	14	867	C	O5'-P-OP1	-7.66	98.80	105.70
26	1H	672	C	OP2-P-O3'	7.66	122.05	105.20
26	1H	1278	A	N1-C6-N6	7.66	123.20	118.60
55	Q8	50	LEU	CA-CB-CG	7.66	132.92	115.30
26	1H	2444	G	C8-N9-C4	-7.65	103.34	106.40
26	14	737	C	N3-C2-O2	7.65	127.26	121.90
26	14	2039	C	C2-N1-C1'	7.65	127.22	118.80
26	1H	2711	A	OP1-P-O3'	7.65	122.03	105.20
26	1H	984	A	O5'-P-OP2	-7.65	98.82	105.70
26	1H	2070	G	C8-N9-C4	7.64	109.46	106.40
26	14	979	G	C8-N9-C4	-7.64	103.34	106.40
1	13	23	C	C6-N1-C2	-7.63	117.25	120.30
27	16	115	G	C5-C6-N1	7.63	115.32	111.50
26	14	2463	C	C6-N1-C2	7.63	123.35	120.30
26	1H	610	C	C5-C6-N1	-7.63	117.19	121.00
26	1H	66	C	C6-N1-C2	-7.63	117.25	120.30
26	1H	1610	A	N9-C4-C5	-7.63	102.75	105.80
26	1H	678	C	N3-C4-C5	7.62	124.95	121.90
26	1H	1299	G	O5'-P-OP1	-7.62	98.84	105.70
1	13	656	C	C5-C6-N1	7.62	124.81	121.00
26	1H	778	G	C5-C6-O6	7.62	133.17	128.60
26	1H	203	C	O5'-P-OP2	7.61	119.84	110.70
26	14	943	U	O5'-P-OP1	-7.61	98.85	105.70
26	1H	265	A	N1-C2-N3	7.61	133.10	129.30
26	1H	1558	A	C2-N3-C4	-7.61	106.80	110.60
26	1H	1817	G	N3-C2-N2	7.61	125.23	119.90
26	1H	1501	C	O5'-P-OP2	7.61	119.83	110.70
26	1H	815	C	C6-N1-C2	7.61	123.34	120.30
1	1G	1499	A	O5'-P-OP1	-7.61	98.85	105.70
26	1H	1332	G	C6-C5-N7	-7.60	125.84	130.40
26	1H	1640	C	O5'-P-OP1	7.60	119.82	110.70
26	1H	1814	G	C4-C5-N7	7.60	113.84	110.80
1	1G	1297	C	P-O3'-C3'	7.60	128.82	119.70
1	1G	1502	A	N7-C8-N9	7.60	117.60	113.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	697	C	O5'-P-OP1	-7.60	98.86	105.70
26	1H	613	U	N1-C2-N3	7.59	119.46	114.90
26	1H	122	G	OP1-P-OP2	7.59	130.99	119.60
1	1G	337	C	C5-C6-N1	7.59	124.80	121.00
26	1H	1210	A	N7-C8-N9	7.59	117.59	113.80
1	13	1517	G	O5'-P-OP2	-7.59	98.87	105.70
24	3L	76	A	O4'-C1'-N9	7.59	114.27	108.20
26	1H	763	G	C6-C5-N7	-7.58	125.85	130.40
26	14	196	A	O4'-C1'-N9	7.58	114.27	108.20
26	1H	1939	U	N3-C2-O2	7.58	127.51	122.20
26	14	201	C	C5-C6-N1	-7.58	117.21	121.00
26	14	776	G	N3-C2-N2	-7.58	114.60	119.90
26	1H	1626	G	N7-C8-N9	7.57	116.89	113.10
26	14	2575	C	C5-C4-N4	7.57	125.50	120.20
26	1H	1265	A	O5'-P-OP1	-7.57	98.89	105.70
26	1H	2420	C	O5'-P-OP1	-7.57	98.89	105.70
26	1H	1397	U	N1-C2-O2	7.57	128.10	122.80
26	1H	827	U	O5'-P-OP1	7.57	119.78	110.70
26	1H	826	U	N1-C2-O2	-7.57	117.50	122.80
1	13	1354	C	C6-N1-C2	-7.56	117.28	120.30
26	1H	478	A	C6-N1-C2	-7.56	114.06	118.60
26	14	2273	A	O5'-P-OP2	-7.56	98.90	105.70
1	13	740	U	O5'-P-OP2	-7.56	98.90	105.70
23	2K	21	U	C2-N3-C4	-7.56	122.47	127.00
26	1H	1382	G	C4-C5-N7	7.55	113.82	110.80
1	13	1158	C	N3-C2-O2	-7.55	116.61	121.90
26	1H	2252	G	N7-C8-N9	-7.55	109.32	113.10
26	14	2255	G	O5'-P-OP2	-7.55	98.90	105.70
1	13	792	A	N7-C8-N9	7.55	117.57	113.80
55	Q8	52	LYS	C-N-CD	-7.55	104.00	120.60
26	1H	1430	C	OP1-P-O3'	7.54	121.79	105.20
26	1H	965	C	C5-C6-N1	7.54	124.77	121.00
26	1H	1621	U	N3-C4-O4	7.54	124.68	119.40
26	1H	1939	U	N1-C2-O2	-7.54	117.52	122.80
26	1H	2275	C	O4'-C1'-N1	-7.53	102.17	108.20
26	1H	2503	A	C2-N3-C4	7.53	114.36	110.60
26	14	2511	U	O5'-P-OP2	-7.53	98.92	105.70
26	14	2609	U	O5'-P-OP2	-7.53	98.92	105.70
26	1H	2330	G	C4-C5-N7	7.52	113.81	110.80
24	3K	71	G	O4'-C1'-N9	7.52	114.22	108.20
26	1H	2424	C	OP1-P-OP2	7.52	130.88	119.60
26	14	2818	G	C5-C6-O6	-7.52	124.09	128.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	1391	U	N3-C4-O4	7.52	124.66	119.40
26	1H	271(B)	G	N3-C4-C5	-7.51	124.84	128.60
26	1H	526	A	N9-C4-C5	7.51	108.81	105.80
26	1H	735	A	N7-C8-N9	-7.51	110.05	113.80
26	1H	137	C	C6-N1-C2	-7.51	117.30	120.30
26	14	2587	A	N1-C6-N6	7.50	123.10	118.60
26	1H	1781	C	N1-C2-O2	7.49	123.40	118.90
26	14	388	G	C5-C6-N1	-7.49	107.75	111.50
26	1H	1546	C	C6-N1-C2	-7.49	117.31	120.30
26	14	1585	C	N1-C2-O2	7.49	123.39	118.90
26	14	1678	G	C4-C5-N7	7.49	113.80	110.80
26	1H	1698	A	N3-C4-C5	7.48	132.04	126.80
26	1H	596	G	N1-C6-O6	7.48	124.39	119.90
26	14	1024	G	N1-C6-O6	7.48	124.39	119.90
26	14	1274	A	C5-C6-N6	-7.48	117.72	123.70
26	14	2392	A	C8-N9-C4	-7.48	102.81	105.80
26	1H	74	A	N7-C8-N9	7.48	117.54	113.80
26	1H	617	G	N7-C8-N9	-7.48	109.36	113.10
26	1H	2586	C	N3-C4-N4	7.48	123.23	118.00
26	14	829	A	O5'-P-OP2	-7.47	98.97	105.70
26	14	2217	G	N1-C6-O6	7.47	124.39	119.90
26	1H	1323	U	OP1-P-OP2	-7.47	108.39	119.60
26	1H	1604	C	O5'-P-OP2	7.47	119.67	110.70
26	1H	528	A	C5-N7-C8	-7.47	100.17	103.90
26	1H	628	G	C8-N9-C4	7.47	109.39	106.40
26	14	530	G	C2-N3-C4	-7.47	108.17	111.90
26	14	1678	G	C2-N3-C4	-7.46	108.17	111.90
26	1H	852	G	O5'-P-OP2	-7.46	98.98	105.70
26	14	2038	G	C8-N9-C4	7.46	109.39	106.40
1	1G	1286	A	C8-N9-C4	-7.46	102.82	105.80
26	1H	1936	A	C5-C6-N6	-7.46	117.73	123.70
26	1H	2830	G	C8-N9-C4	-7.46	103.42	106.40
26	1H	2275	C	OP1-P-O3'	7.46	121.60	105.20
36	35	65	ARG	N-CA-C	-7.45	90.88	111.00
26	1H	446	G	N1-C6-O6	7.45	124.37	119.90
26	1H	271(B)	G	P-O3'-C3'	7.45	128.64	119.70
26	1H	859	G	N3-C4-C5	7.45	132.32	128.60
26	14	676	A	N1-C6-N6	7.45	123.07	118.60
26	14	1332	G	O4'-C1'-N9	-7.45	102.24	108.20
26	1H	252	G	O5'-P-OP1	7.44	119.63	110.70
26	1H	1128	A	O5'-P-OP2	-7.44	99.00	105.70
26	1H	1307	A	N1-C6-N6	7.44	123.06	118.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	1G	1484	C	O5'-P-OP2	-7.44	99.00	105.70
23	2L	29	C	C5-C6-N1	7.44	124.72	121.00
26	14	1256	G	N1-C6-O6	7.44	124.36	119.90
26	14	2313	C	C6-N1-C2	-7.44	117.32	120.30
26	1H	1027	A	C2-N3-C4	-7.44	106.88	110.60
26	14	770	G	OP1-P-OP2	-7.44	108.44	119.60
26	14	2351	G	N3-C4-N9	7.44	130.46	126.00
26	1H	2401	U	C5-C6-N1	7.44	126.42	122.70
1	1G	183	G	N1-C6-O6	7.44	124.36	119.90
26	14	1274	A	C6-C5-N7	-7.43	127.10	132.30
26	14	1325	G	N3-C4-N9	7.43	130.46	126.00
26	1H	119	A	C4-C5-N7	-7.43	106.98	110.70
26	1H	946	G	N7-C8-N9	-7.43	109.38	113.10
26	1H	826	U	C4-C5-C6	7.43	124.16	119.70
26	1H	1123	C	N1-C2-O2	-7.43	114.44	118.90
26	14	684	G	C8-N9-C4	-7.43	103.43	106.40
26	14	1256	G	C5-C6-O6	-7.43	124.14	128.60
26	1H	223	A	O5'-P-OP2	-7.42	99.02	105.70
26	1H	1806	C	OP1-P-OP2	7.42	130.73	119.60
26	1H	1691	C	C6-N1-C2	-7.42	117.33	120.30
26	1H	140	A	C2-N3-C4	-7.42	106.89	110.60
1	13	295	C	O5'-P-OP2	-7.42	99.03	105.70
1	13	329	A	O5'-P-OP2	-7.41	99.03	105.70
26	1H	462	C	O5'-P-OP2	-7.41	99.03	105.70
26	1H	464	U	C5-C6-N1	-7.41	119.00	122.70
26	1H	860	U	C6-N1-C1'	-7.41	110.83	121.20
1	13	789	U	N1-C2-N3	7.41	119.34	114.90
26	14	1630(A)	C	N1-C2-O2	-7.41	114.46	118.90
1	13	687	A	P-O3'-C3'	7.40	128.58	119.70
26	1H	190	A	C6-N1-C2	-7.40	114.16	118.60
1	1G	320	C	C6-N1-C2	7.40	123.26	120.30
26	14	330	A	N1-C6-N6	7.40	123.04	118.60
26	1H	1914	C	C6-N1-C2	-7.40	117.34	120.30
26	14	2439	A	N1-C6-N6	7.40	123.04	118.60
26	1H	1940	U	N3-C4-O4	7.40	124.58	119.40
26	14	2301	C	C6-N1-C2	-7.39	117.34	120.30
26	1H	529	A	C8-N9-C4	-7.39	102.84	105.80
34	15	42	TRP	NE1-CE2-CD2	7.39	114.69	107.30
26	1H	2585	U	N3-C2-O2	-7.38	117.03	122.20
26	1H	1376	C	C6-N1-C2	-7.38	117.35	120.30
26	1H	2822	G	C6-C5-N7	-7.38	125.97	130.40
1	13	690	G	N3-C4-N9	7.38	130.43	126.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	1605	C	C2-N3-C4	-7.38	116.21	119.90
26	1H	1367	A	C2-N3-C4	-7.38	106.91	110.60
1	13	535	A	N1-C6-N6	-7.38	114.17	118.60
26	1H	1443	G	C6-C5-N7	-7.38	125.97	130.40
26	14	113	G	N1-C6-O6	7.37	124.33	119.90
26	1H	1792	G	N1-C6-O6	-7.37	115.48	119.90
26	1H	2395	C	C2-N1-C1'	7.37	126.90	118.80
1	1G	1518	A	O5'-P-OP1	-7.37	99.07	105.70
26	14	113	G	N3-C4-C5	7.37	132.28	128.60
26	1H	663	G	C4-C5-C6	7.37	123.22	118.80
26	1H	2619	C	O5'-P-OP2	-7.37	99.07	105.70
26	14	1270	C	C6-N1-C2	-7.36	117.35	120.30
26	14	2001	A	C5-C6-N6	-7.36	117.81	123.70
26	14	1332	G	C4-C5-C6	7.36	123.22	118.80
26	14	2332	U	O5'-P-OP1	7.36	119.53	110.70
26	1H	763	G	C5-C6-O6	-7.36	124.18	128.60
26	14	1939	U	OP2-P-O3'	7.36	121.39	105.20
26	14	1616	A	O4'-C1'-N9	7.36	114.09	108.20
26	1H	37	C	N3-C4-C5	-7.36	118.96	121.90
26	14	1377	G	C8-N9-C4	-7.36	103.46	106.40
26	1H	1831	G	O5'-P-OP2	7.36	119.53	110.70
26	1H	811	U	O5'-P-OP1	-7.35	99.08	105.70
26	1H	762	U	N1-C2-O2	7.35	127.94	122.80
26	14	1742	C	C6-N1-C2	-7.35	117.36	120.30
1	13	219	C	C6-N1-C2	-7.35	117.36	120.30
26	14	74	A	C5-N7-C8	-7.35	100.23	103.90
26	1H	698	C	C4-C5-C6	7.34	121.07	117.40
26	14	2513	G	C5-C6-O6	-7.34	124.19	128.60
26	14	1313	U	C6-N1-C2	-7.34	116.59	121.00
23	2K	35	C	C2-N1-C1'	7.34	126.87	118.80
26	1H	783	A	N1-C6-N6	7.34	123.00	118.60
26	1H	1229(A)	G	O5'-P-OP2	-7.34	99.09	105.70
26	1H	537	C	O5'-P-OP2	-7.34	99.09	105.70
1	13	1199	U	N3-C4-C5	-7.34	110.20	114.60
26	1H	2446	G	C5-C6-O6	-7.34	124.20	128.60
26	14	1527	G	N3-C4-C5	7.33	132.27	128.60
26	1H	113	G	N1-C6-O6	7.33	124.30	119.90
26	1H	1817	G	N1-C2-N2	-7.33	109.60	116.20
26	14	528	A	N1-C6-N6	7.33	123.00	118.60
26	1H	762	U	C2-N1-C1'	7.33	126.49	117.70
26	1H	1521	G	C8-N9-C4	-7.33	103.47	106.40
26	1H	2080	G	O5'-P-OP1	-7.33	99.11	105.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	1G	27	G	N1-C6-O6	7.33	124.30	119.90
26	1H	1967	C	N1-C2-O2	7.33	123.30	118.90
26	14	2346	A	C2-N3-C4	-7.33	106.94	110.60
27	16	81	G	C4-C5-N7	7.33	113.73	110.80
26	1H	967	C	O5'-P-OP2	-7.32	99.11	105.70
26	14	2087	G	N9-C4-C5	-7.32	102.47	105.40
26	1H	671	C	C6-N1-C2	7.32	123.23	120.30
26	14	945	A	N9-C4-C5	-7.32	102.87	105.80
26	14	774	A	C5-N7-C8	-7.31	100.24	103.90
26	14	2249	U	N3-C4-C5	-7.31	110.21	114.60
26	1H	2499	C	N3-C2-O2	7.31	127.02	121.90
26	1H	787	U	OP1-P-OP2	-7.31	108.64	119.60
26	1H	2060	A	N1-C6-N6	-7.31	114.22	118.60
26	1H	2379	G	N3-C4-N9	7.31	130.38	126.00
26	1H	1369	G	N3-C4-C5	-7.30	124.95	128.60
26	1H	2546	U	N1-C2-O2	-7.30	117.69	122.80
26	1H	1308	A	N1-C2-N3	7.30	132.95	129.30
26	1H	559	G	N1-C6-O6	7.30	124.28	119.90
26	1H	845	G	P-O3'-C3'	7.30	128.46	119.70
26	1H	2276	G	N3-C2-N2	-7.30	114.79	119.90
26	14	1284	A	O5'-P-OP2	-7.30	99.13	105.70
26	1H	446	G	N3-C4-N9	7.30	130.38	126.00
26	1H	1772	G	N3-C2-N2	7.29	125.01	119.90
26	1H	1681	G	N1-C6-O6	7.29	124.28	119.90
26	14	2386	C	C6-N1-C2	7.29	123.22	120.30
26	14	922	U	C5-C6-N1	7.29	126.34	122.70
26	1H	974(A)	C	N3-C2-O2	-7.29	116.80	121.90
26	1H	1691	C	OP1-P-O3'	7.29	121.23	105.20
1	1G	995	C	C6-N1-C2	-7.29	117.39	120.30
26	14	2301	C	C5-C6-N1	7.29	124.64	121.00
1	13	115	G	C8-N9-C4	-7.28	103.49	106.40
26	14	2776	A	C8-N9-C4	-7.28	102.89	105.80
26	1H	2586	C	C5-C4-N4	-7.28	115.10	120.20
26	14	2507	C	O5'-P-OP2	-7.28	99.15	105.70
1	1G	528	C	O4'-C1'-N1	7.28	114.02	108.20
26	1H	2497	A	C6-N1-C2	-7.28	114.23	118.60
26	1H	1854	A	N1-C6-N6	-7.28	114.23	118.60
26	1H	2582	G	C6-C5-N7	-7.28	126.03	130.40
26	14	2012	G	N1-C6-O6	7.28	124.27	119.90
26	1H	2415	G	N1-C6-O6	7.27	124.26	119.90
28	19	37	LEU	CA-CB-CG	7.27	132.03	115.30
26	1H	1332	G	N1-C6-O6	7.27	124.26	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	774	A	C8-N9-C1'	7.27	140.79	127.70
1	13	1505	G	OP1-P-OP2	-7.27	108.70	119.60
26	1H	1820	U	C6-N1-C2	7.27	125.36	121.00
26	14	113	G	C5-C6-O6	-7.27	124.24	128.60
26	14	1790	C	C6-N1-C2	7.27	123.21	120.30
1	13	1335	C	C6-N1-C2	7.26	123.20	120.30
26	1H	1410	G	C8-N9-C4	7.26	109.30	106.40
26	1H	1698	A	C4-C5-N7	7.26	114.33	110.70
1	13	280	C	C6-N1-C2	7.26	123.20	120.30
26	1H	107	C	C6-N1-C2	7.25	123.20	120.30
26	1H	729	G	N7-C8-N9	7.25	116.73	113.10
26	1H	688	U	N1-C2-N3	7.25	119.25	114.90
26	1H	1380	G	N1-C6-O6	7.25	124.25	119.90
26	14	2080	G	C5-C6-O6	7.25	132.95	128.60
26	1H	1428	C	N3-C4-C5	7.25	124.80	121.90
26	14	2444	G	C5-C6-O6	7.25	132.95	128.60
26	1H	146	G	N1-C6-O6	7.25	124.25	119.90
26	14	37	C	C6-N1-C2	-7.24	117.40	120.30
26	14	1681	G	N7-C8-N9	7.24	116.72	113.10
26	1H	470	A	N1-C6-N6	7.24	122.94	118.60
26	14	2080	G	C4-C5-N7	-7.24	107.91	110.80
1	1G	690	G	C5-N7-C8	-7.24	100.68	104.30
26	14	2623	G	O5'-P-OP1	7.24	119.38	110.70
26	14	2062	A	N1-C6-N6	7.23	122.94	118.60
26	1H	965	C	C6-N1-C2	-7.23	117.41	120.30
26	14	1653	G	O5'-P-OP2	-7.23	99.19	105.70
27	16	56	G	C8-N9-C4	-7.23	103.51	106.40
26	14	499	U	N3-C2-O2	-7.23	117.14	122.20
26	1H	2004	G	O5'-P-OP2	-7.22	99.20	105.70
26	1H	2847	U	O5'-P-OP1	-7.22	99.20	105.70
26	14	729	G	N1-C6-O6	7.22	124.23	119.90
26	14	2364	C	O5'-P-OP2	-7.22	99.20	105.70
26	14	1821	A	N1-C6-N6	7.22	122.93	118.60
26	1H	2057	A	C8-N9-C4	7.22	108.69	105.80
26	1H	663	G	N7-C8-N9	7.22	116.71	113.10
26	1H	2413	G	N1-C6-O6	7.21	124.23	119.90
26	14	2206	C	O5'-P-OP2	-7.21	99.21	105.70
26	1H	1271	G	O5'-P-OP2	-7.21	99.21	105.70
26	14	2623	G	N3-C4-C5	-7.21	124.99	128.60
26	14	1971	A	C8-N9-C4	7.21	108.69	105.80
26	14	2873	A	C4-C5-N7	7.21	114.31	110.70
26	1H	2567	G	O5'-P-OP1	-7.21	99.21	105.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	2491	U	O5'-P-OP1	-7.21	99.21	105.70
26	1H	208	C	C5-C6-N1	-7.21	117.40	121.00
26	1H	636	G	O5'-P-OP1	-7.21	99.21	105.70
23	2K	21	U	C4-C5-C6	-7.20	115.38	119.70
26	1H	2406	U	O4'-C1'-N1	-7.20	102.44	108.20
1	1G	284	G	N1-C6-O6	7.20	124.22	119.90
26	1H	2700	C	N3-C4-C5	7.20	124.78	121.90
26	1H	696	G	O5'-P-OP2	7.20	119.34	110.70
24	3K	76	A	C6-C5-N7	-7.19	127.26	132.30
1	13	63	C	C6-N1-C2	-7.19	117.42	120.30
26	1H	1764	G	C8-N9-C4	-7.19	103.52	106.40
26	14	298	G	N1-C6-O6	7.19	124.22	119.90
1	13	956	U	C6-N1-C2	-7.19	116.69	121.00
26	1H	689	A	C2-N3-C4	-7.19	107.00	110.60
26	1H	2011	U	N3-C2-O2	7.19	127.23	122.20
26	14	1429	G	C8-N9-C4	-7.19	103.53	106.40
26	1H	810	U	O5'-P-OP2	-7.18	99.23	105.70
26	1H	822	U	N3-C2-O2	-7.18	117.17	122.20
26	1H	1653	G	O5'-P-OP2	-7.18	99.23	105.70
26	14	2607	G	N9-C4-C5	-7.18	102.53	105.40
26	1H	198	C	N1-C2-O2	7.18	123.21	118.90
24	3K	76	A	N1-C6-N6	7.18	122.91	118.60
26	1H	587	C	O5'-P-OP1	-7.18	99.24	105.70
26	1H	1279	G	O5'-P-OP1	7.18	119.31	110.70
26	1H	1622	G	C4-C5-N7	-7.18	107.93	110.80
26	1H	1772	G	N1-C2-N2	-7.18	109.74	116.20
26	1H	446	G	C8-N9-C1'	-7.17	117.67	127.00
26	1H	691	C	C5-C6-N1	-7.17	117.41	121.00
26	1H	2274	A	OP2-P-O3'	7.17	120.98	105.20
26	14	1336	A	O5'-P-OP2	-7.17	99.25	105.70
26	14	2038	G	N9-C4-C5	-7.17	102.53	105.40
1	13	5	U	P-O3'-C3'	7.17	128.31	119.70
1	13	122	G	N1-C6-O6	7.17	124.20	119.90
26	14	252	G	O5'-P-OP2	-7.17	99.25	105.70
26	14	1933	G	O5'-P-OP2	-7.17	99.25	105.70
26	1H	2506	U	N1-C2-O2	7.17	127.82	122.80
26	14	1239	G	O5'-P-OP1	-7.17	99.25	105.70
26	14	1782	C	C5-C4-N4	-7.16	115.19	120.20
26	1H	1786	A	OP1-P-O3'	7.16	120.95	105.20
26	1H	330	A	C5-N7-C8	-7.16	100.32	103.90
26	1H	2238	G	OP1-P-OP2	7.16	130.34	119.60
26	1H	1300	U	O5'-P-OP2	-7.16	99.26	105.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	2859	G	C8-N9-C4	-7.16	103.54	106.40
26	1H	1611	C	C2-N3-C4	-7.15	116.32	119.90
26	14	1640	C	O5'-P-OP1	7.15	119.28	110.70
1	13	1199	U	C5-C4-O4	7.15	130.19	125.90
26	1H	839	U	N3-C4-C5	-7.15	110.31	114.60
26	14	752	A	C8-N9-C4	-7.15	102.94	105.80
26	14	2072	G	OP1-P-OP2	-7.15	108.87	119.60
26	1H	446	G	C8-N9-C4	7.15	109.26	106.40
26	14	801	G	C5-C6-O6	7.15	132.89	128.60
26	14	1698	A	C5-C6-N6	-7.15	117.98	123.70
26	14	1698	A	C4-C5-C6	7.15	120.57	117.00
26	1H	983	A	C8-N9-C4	7.14	108.66	105.80
27	16	99	A	OP1-P-OP2	7.14	130.31	119.60
26	1H	1332	G	N1-C2-N3	7.13	128.18	123.90
1	1G	529	G	C5-C6-O6	-7.13	124.32	128.60
26	14	1962	C	N3-C4-C5	7.13	124.75	121.90
26	14	2087	G	C8-N9-C4	7.13	109.25	106.40
26	1H	2073	C	OP2-P-O3'	7.13	120.89	105.20
26	1H	2847	U	O5'-P-OP2	7.13	119.26	110.70
23	2L	35	C	C2-N1-C1'	7.13	126.64	118.80
1	13	1266	G	N3-C4-C5	7.13	132.16	128.60
26	1H	299	A	OP2-P-O3'	7.12	120.88	105.20
26	1H	704	G	N9-C4-C5	7.12	108.25	105.40
26	1H	1800	C	O5'-P-OP2	7.12	119.25	110.70
1	13	564	C	N3-C4-C5	-7.12	119.05	121.90
26	1H	1998	G	C8-N9-C4	7.12	109.25	106.40
26	1H	2681	C	N3-C2-O2	-7.11	116.92	121.90
26	1H	845	G	N3-C4-C5	7.10	132.15	128.60
26	1H	954	G	N3-C2-N2	-7.10	114.93	119.90
26	1H	2318	G	O4'-C1'-N9	7.10	113.88	108.20
26	14	2001	A	N9-C4-C5	-7.10	102.96	105.80
26	14	783	A	O5'-P-OP2	-7.10	99.31	105.70
26	1H	575	A	O5'-P-OP1	-7.10	99.31	105.70
26	1H	2367	G	C8-N9-C4	-7.09	103.56	106.40
26	14	2452	C	N3-C2-O2	7.09	126.86	121.90
26	14	1313	U	O5'-P-OP2	-7.09	99.32	105.70
26	14	1964	G	O5'-P-OP1	-7.09	99.32	105.70
26	14	855	G	C8-N9-C4	-7.09	103.56	106.40
26	1H	974(A)	C	C2-N3-C4	7.09	123.44	119.90
26	1H	2314	C	O5'-P-OP1	7.09	119.20	110.70
26	1H	2697	G	OP1-P-OP2	7.09	130.23	119.60
25	4K	18	G	C8-N9-C4	-7.08	103.57	106.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	2492	U	N1-C2-O2	7.08	127.76	122.80
26	1H	842	G	C5-C6-O6	-7.08	124.35	128.60
26	1H	1573	G	OP2-P-O3'	7.08	120.78	105.20
1	13	1498	U	C2-N1-C1'	7.08	126.19	117.70
26	14	912	C	C6-N1-C2	-7.08	117.47	120.30
26	1H	2249	U	C6-N1-C2	-7.08	116.75	121.00
26	1H	788	A	C8-N9-C4	7.07	108.63	105.80
26	1H	956	G	N1-C6-O6	7.07	124.14	119.90
26	14	1934	C	N1-C2-O2	7.07	123.14	118.90
26	1H	2638	G	N3-C4-N9	7.07	130.24	126.00
26	14	510	C	O5'-P-OP2	-7.07	99.34	105.70
26	14	2688	U	N1-C2-N3	7.07	119.14	114.90
26	1H	1274	A	C8-N9-C4	-7.07	102.97	105.80
26	1H	1787	A	O4'-C1'-N9	-7.07	102.55	108.20
26	1H	1825	A	C5-C6-N6	7.07	129.35	123.70
26	1H	189	G	N7-C8-N9	-7.06	109.57	113.10
26	1H	2346	A	C4-C5-C6	7.06	120.53	117.00
26	14	728	G	N3-C4-N9	7.06	130.24	126.00
26	14	955	C	C6-N1-C2	-7.06	117.47	120.30
26	14	2439	A	P-O3'-C3'	7.06	128.18	119.70
26	14	570	G	C4-N9-C1'	7.06	135.68	126.50
26	1H	1777	U	C4-C5-C6	7.06	123.94	119.70
26	1H	74	A	N3-C4-C5	7.06	131.74	126.80
26	1H	778	G	O5'-P-OP1	7.06	119.17	110.70
26	1H	1443	G	C5-C6-N1	-7.06	107.97	111.50
26	14	2256	G	O5'-P-OP2	-7.05	99.35	105.70
26	14	1347	G	OP1-P-O3'	7.05	120.71	105.20
26	1H	1958	C	OP1-P-O3'	7.05	120.70	105.20
26	1H	2048	G	C8-N9-C4	-7.05	103.58	106.40
26	14	1380	G	N1-C6-O6	7.05	124.13	119.90
26	1H	71	A	C6-C5-N7	-7.04	127.37	132.30
26	14	676	A	N3-C4-N9	-7.04	121.77	127.40
26	14	2526	G	N3-C4-N9	-7.04	121.78	126.00
26	1H	966	G	N1-C6-O6	-7.04	115.68	119.90
26	1H	774	A	N9-C4-C5	7.04	108.61	105.80
26	14	130	C	N3-C4-C5	7.04	124.72	121.90
26	1H	49	A	N7-C8-N9	-7.04	110.28	113.80
26	14	1608	A	N1-C6-N6	-7.04	114.38	118.60
1	13	186(A)	C	C6-N1-C2	-7.04	117.49	120.30
26	1H	1428	C	C5-C6-N1	-7.04	117.48	121.00
26	14	573	G	C2-N3-C4	7.03	115.42	111.90
26	14	1827	C	N3-C2-O2	-7.03	116.98	121.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	583	G	C8-N9-C4	-7.03	103.59	106.40
26	1H	1569	A	C8-N9-C4	-7.03	102.99	105.80
26	1H	2502	G	N3-C4-C5	-7.03	125.09	128.60
26	1H	2607	G	C2-N3-C4	-7.03	108.39	111.90
26	14	2374	C	C6-N1-C2	7.02	123.11	120.30
26	1H	321	G	N3-C4-N9	7.02	130.21	126.00
26	1H	1858	G	N1-C6-O6	7.02	124.11	119.90
26	1H	2428	G	C8-N9-C4	-7.02	103.59	106.40
1	1G	354	G	C6-C5-N7	-7.02	126.19	130.40
1	13	738	C	C6-N1-C2	-7.02	117.49	120.30
1	13	966	G	C8-N9-C4	7.01	109.21	106.40
26	1H	753	C	C6-N1-C2	-7.01	117.49	120.30
26	1H	1296	G	OP2-P-O3'	7.01	120.63	105.20
26	1H	1610	A	N1-C6-N6	7.01	122.81	118.60
26	1H	1698	A	C5-C6-N1	-7.01	114.19	117.70
1	13	872	A	O4'-C1'-N9	7.01	113.81	108.20
26	1H	1325	G	C6-C5-N7	-7.01	126.19	130.40
26	14	209	C	N3-C4-N4	7.01	122.91	118.00
26	1H	624	C	N1-C2-O2	-7.01	114.69	118.90
26	1H	1971	A	C6-N1-C2	-7.01	114.39	118.60
26	1H	2747	G	N9-C4-C5	-7.01	102.60	105.40
26	1H	679	C	O5'-P-OP1	-7.01	99.39	105.70
26	1H	1305	C	N1-C2-O2	7.00	123.10	118.90
26	14	825	C	N3-C4-C5	-7.00	119.10	121.90
26	14	2324	C	N3-C4-C5	7.00	124.70	121.90
26	1H	1363	C	N3-C4-C5	7.00	124.70	121.90
1	1G	1354	C	C6-N1-C2	-7.00	117.50	120.30
26	14	2386	C	C5-C6-N1	-7.00	117.50	121.00
26	14	2512	C	C6-N1-C2	7.00	123.10	120.30
26	1H	732	C	N1-C2-O2	-7.00	114.70	118.90
26	1H	1574	C	OP2-P-O3'	7.00	120.59	105.20
26	1H	664	C	C2-N3-C4	-7.00	116.40	119.90
26	1H	2609	U	C5-C6-N1	-6.99	119.20	122.70
1	13	525	C	N3-C4-N4	6.99	122.89	118.00
26	1H	1614	A	C5-N7-C8	-6.99	100.41	103.90
26	1H	2712	U	C2-N3-C4	-6.99	122.81	127.00
26	1H	2499	C	C5-C4-N4	-6.99	115.31	120.20
26	14	602	G	N1-C6-O6	6.99	124.09	119.90
26	14	1258	C	N3-C2-O2	6.99	126.79	121.90
26	1H	1204	A	C6-C5-N7	-6.99	127.41	132.30
26	1H	2449	U	C6-N1-C2	-6.99	116.81	121.00
1	1G	567	G	N1-C6-O6	-6.99	115.71	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	13	1519	A	C5-C6-N1	-6.98	114.21	117.70
26	1H	1971	A	C2-N3-C4	6.98	114.09	110.60
26	14	510	C	C6-N1-C2	-6.98	117.51	120.30
26	14	2518	A	C5-C6-N6	-6.98	118.12	123.70
26	14	1786	A	C5-C6-N1	-6.98	114.21	117.70
26	14	829	A	O5'-P-OP1	-6.98	99.42	105.70
26	1H	144	C	C2-N3-C4	-6.97	116.41	119.90
26	1H	733	G	N1-C2-N2	-6.97	109.92	116.20
23	2L	40	C	C6-N1-C2	-6.97	117.51	120.30
1	13	266	G	N1-C6-O6	6.97	124.08	119.90
26	1H	1537	C	C6-N1-C2	-6.97	117.51	120.30
26	1H	818	G	O5'-P-OP1	-6.97	99.43	105.70
26	1H	134	C	N3-C2-O2	-6.96	117.03	121.90
26	1H	1298	C	OP1-P-O3'	6.96	120.52	105.20
26	1H	2616	C	N1-C2-O2	-6.96	114.72	118.90
26	14	1142	U	C2-N1-C1'	6.96	126.06	117.70
26	1H	1278	A	C5-C6-N6	-6.96	118.13	123.70
26	1H	2598	A	OP2-P-O3'	6.96	120.52	105.20
26	1H	1399	C	C6-N1-C2	-6.96	117.52	120.30
23	2K	22	A	O5'-P-OP2	-6.96	99.44	105.70
26	1H	446	G	C6-C5-N7	-6.96	126.22	130.40
26	1H	837	C	C2-N1-C1'	6.96	126.45	118.80
26	14	729	G	C5-C6-O6	-6.96	124.43	128.60
26	1H	1764	G	N9-C4-C5	6.95	108.18	105.40
26	1H	1776	G	N9-C4-C5	-6.95	102.62	105.40
26	1H	2395	C	C6-N1-C1'	-6.95	112.46	120.80
26	14	1899	G	N1-C2-N3	6.95	128.07	123.90
26	1H	1621	U	N1-C2-O2	-6.95	117.94	122.80
26	1H	2575	C	C4-C5-C6	6.95	120.87	117.40
1	1G	317	G	N1-C6-O6	6.95	124.07	119.90
26	14	1902	C	O5'-P-OP2	6.95	119.04	110.70
26	14	2335	A	N1-C6-N6	-6.95	114.43	118.60
26	1H	1129	A	O5'-P-OP2	-6.95	99.45	105.70
26	14	1359	A	N9-C4-C5	-6.95	103.02	105.80
26	1H	689	A	N1-C2-N3	6.94	132.77	129.30
26	1H	925	C	O5'-P-OP2	-6.94	99.45	105.70
1	13	513	C	C5-C6-N1	6.94	124.47	121.00
26	1H	698	C	C5-C6-N1	-6.94	117.53	121.00
26	1H	1308	A	C6-N1-C2	-6.94	114.44	118.60
26	1H	1373	A	C8-N9-C4	6.94	108.58	105.80
26	1H	705	A	O5'-P-OP2	-6.94	99.45	105.70
26	1H	1621	U	N3-C2-O2	6.94	127.06	122.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	13	529	G	C5-C6-O6	-6.93	124.44	128.60
26	1H	1969	A	N1-C6-N6	-6.93	114.44	118.60
26	1H	2584	U	N3-C4-O4	-6.93	114.55	119.40
1	13	1266	G	C4-N9-C1'	-6.93	117.49	126.50
26	1H	2609	U	C4-C5-C6	6.93	123.86	119.70
26	14	971	C	C6-N1-C2	-6.93	117.53	120.30
26	1H	988	A	C8-N9-C4	-6.93	103.03	105.80
26	1H	2726	U	C5-C6-N1	-6.93	119.24	122.70
26	14	819	A	C8-N9-C4	-6.92	103.03	105.80
26	1H	774	A	C4-C5-N7	6.92	114.16	110.70
26	1H	1427	A	N1-C6-N6	-6.92	114.45	118.60
26	1H	691	C	N3-C2-O2	6.92	126.74	121.90
26	1H	1468	C	C6-N1-C2	-6.92	117.53	120.30
26	14	1029	A	N1-C6-N6	6.92	122.75	118.60
26	1H	2292	C	OP1-P-OP2	-6.92	109.22	119.60
26	1H	747	U	O5'-P-OP1	-6.92	99.47	105.70
26	1H	779	U	C5-C4-O4	-6.92	121.75	125.90
26	1H	2705	A	C5-C6-N6	-6.92	118.17	123.70
1	1G	818	G	C4-C5-N7	-6.92	108.03	110.80
26	1H	778	G	C4-C5-N7	-6.92	108.03	110.80
26	14	1314	C	C2-N1-C1'	6.92	126.41	118.80
26	1H	106	C	C6-N1-C2	-6.91	117.53	120.30
26	14	2012	G	C5-C6-O6	-6.91	124.45	128.60
26	14	1204	A	O4'-C1'-N9	6.91	113.73	108.20
26	14	2420	C	O5'-P-OP1	-6.91	99.48	105.70
26	1H	599	G	C8-N9-C4	6.91	109.16	106.40
1	13	1199	U	C6-N1-C2	-6.91	116.86	121.00
1	1G	332	G	C8-N9-C4	6.91	109.16	106.40
26	14	1379	A	N1-C6-N6	6.91	122.74	118.60
26	1H	493	G	C5-C6-N1	-6.91	108.05	111.50
26	1H	1660	C	N3-C4-C5	6.91	124.66	121.90
26	14	2688	U	C4-C5-C6	6.91	123.84	119.70
26	1H	1604	C	C6-N1-C2	-6.90	117.54	120.30
1	1G	111	G	N1-C6-O6	6.90	124.04	119.90
26	1H	146	G	N9-C4-C5	-6.90	102.64	105.40
26	1H	2609	U	C2-N3-C4	-6.90	122.86	127.00
26	14	1671	U	OP1-P-OP2	6.90	129.95	119.60
37	45	82	ARG	N-CA-C	6.90	129.63	111.00
26	14	252	G	C2-N3-C4	6.90	115.35	111.90
26	14	575	A	O5'-P-OP1	-6.90	99.49	105.70
26	1H	670	A	O4'-C1'-N9	-6.90	102.68	108.20
26	1H	1202	C	N1-C2-O2	-6.90	114.76	118.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	2403	C	N1-C2-O2	-6.90	114.76	118.90
26	1H	429	A	N1-C2-N3	6.89	132.75	129.30
26	14	676	A	OP1-P-OP2	6.89	129.94	119.60
26	1H	34	C	N1-C2-O2	6.89	123.04	118.90
26	1H	2757	A	O5'-P-OP2	-6.89	99.50	105.70
26	1H	456	C	O5'-P-OP2	-6.89	99.50	105.70
26	1H	2486	G	O5'-P-OP2	-6.89	99.50	105.70
26	1H	1248	G	N3-C4-C5	6.89	132.04	128.60
26	14	388	G	N3-C4-N9	-6.89	121.87	126.00
26	1H	820	A	N1-C6-N6	-6.89	114.47	118.60
26	1H	133	C	C5-C6-N1	-6.89	117.56	121.00
26	14	1826	G	C8-N9-C4	6.89	109.15	106.40
26	1H	703	U	C5-C4-O4	6.88	130.03	125.90
26	1H	1657	C	C6-N1-C2	-6.88	117.55	120.30
1	1G	900	A	O5'-P-OP1	-6.88	99.51	105.70
26	14	602	G	C8-N9-C1'	-6.88	118.05	127.00
26	14	2272	U	N1-C2-O2	6.88	127.62	122.80
1	13	1224	G	O5'-P-OP1	6.88	118.96	110.70
26	1H	672	C	O5'-P-OP1	6.88	118.96	110.70
26	1H	2597	G	C4-C5-N7	6.88	113.55	110.80
1	13	101	A	N7-C8-N9	6.88	117.24	113.80
26	1H	1520	U	C6-N1-C2	-6.88	116.87	121.00
26	14	1899	G	C5-C6-O6	6.88	132.73	128.60
26	14	2315	G	N3-C4-N9	6.88	130.12	126.00
26	14	2447	G	P-O3'-C3'	6.87	127.95	119.70
26	1H	632	A	O5'-P-OP2	6.87	118.94	110.70
26	1H	1805	U	N3-C4-O4	6.87	124.21	119.40
26	1H	2418	A	OP1-P-OP2	-6.87	109.30	119.60
26	1H	2518	A	C6-C5-N7	-6.87	127.49	132.30
26	1H	1632	A	C4-C5-N7	6.87	114.13	110.70
26	1H	1914	C	N3-C2-O2	-6.87	117.09	121.90
26	1H	2439	A	N1-C6-N6	6.87	122.72	118.60
23	2K	58	A	O5'-P-OP2	6.86	118.94	110.70
26	1H	826	U	N1-C2-N3	6.86	119.02	114.90
26	14	1598	C	O5'-P-OP2	6.86	118.94	110.70
26	1H	239	U	C5-C4-O4	6.86	130.02	125.90
26	1H	1534	G	C8-N9-C4	-6.86	103.66	106.40
26	1H	1764	G	C5-C6-O6	6.86	132.72	128.60
26	14	204	A	C6-N1-C2	-6.86	114.48	118.60
26	14	2073	C	N1-C2-O2	-6.86	114.78	118.90
26	1H	621	A	O4'-C1'-N9	6.86	113.69	108.20
35	68	22	ILE	CG1-CB-CG2	-6.86	96.31	111.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	1942	C	C5-C6-N1	6.86	124.43	121.00
1	13	535	A	N9-C4-C5	6.85	108.54	105.80
26	14	775	G	N3-C4-C5	-6.85	125.17	128.60
27	16	5	C	C6-N1-C2	6.85	123.04	120.30
22	1K	38	A	N1-C6-N6	6.85	122.71	118.60
23	2K	6	G	C8-N9-C4	6.85	109.14	106.40
26	1H	673	C	OP1-P-OP2	-6.85	109.33	119.60
26	1H	687	C	O5'-P-OP1	-6.85	99.54	105.70
26	14	1298	C	N1-C2-O2	6.85	123.01	118.90
26	1H	621	A	C5-C6-N6	-6.84	118.22	123.70
1	13	1475	G	O5'-P-OP1	-6.84	99.54	105.70
26	1H	594	U	C5-C6-N1	-6.84	119.28	122.70
26	14	1782	C	N3-C4-N4	6.84	122.79	118.00
1	13	30	U	N3-C2-O2	6.84	126.98	122.20
26	1H	1195	G	N3-C2-N2	-6.83	115.12	119.90
26	1H	2199	A	N1-C6-N6	-6.83	114.50	118.60
26	1H	2264	C	OP1-P-O3'	6.83	120.23	105.20
26	1H	2552	U	N1-C2-O2	-6.83	118.02	122.80
26	1H	2271	G	N3-C4-N9	6.82	130.09	126.00
26	14	1332	G	OP1-P-O3'	6.82	120.21	105.20
26	1H	1611	C	O5'-P-OP2	6.82	118.88	110.70
27	16	81	G	C5-N7-C8	-6.82	100.89	104.30
26	1H	2062	A	N9-C4-C5	-6.82	103.07	105.80
26	1H	621	A	N3-C4-C5	6.82	131.57	126.80
26	1H	667	U	N3-C4-O4	6.82	124.17	119.40
26	1H	2031	A	C2-N3-C4	6.82	114.01	110.60
26	1H	1888	G	N3-C4-N9	6.81	130.09	126.00
26	14	1154	G	C5-C6-O6	-6.81	124.51	128.60
26	1H	1604	C	O5'-P-OP1	-6.81	99.57	105.70
26	1H	1830	C	N1-C2-O2	-6.81	114.81	118.90
26	1H	913	U	N3-C4-C5	6.81	118.69	114.60
26	1H	2258	C	N1-C2-O2	-6.81	114.82	118.90
26	1H	2508	G	C8-N9-C4	-6.81	103.68	106.40
26	14	2513	G	N1-C6-O6	6.81	123.98	119.90
1	13	560	U	C5-C6-N1	6.80	126.10	122.70
26	1H	766	C	N1-C2-O2	-6.80	114.82	118.90
26	1H	1265	A	C8-N9-C4	-6.80	103.08	105.80
26	1H	1678	G	N1-C2-N3	6.80	127.98	123.90
1	1G	1502	A	C8-N9-C4	-6.80	103.08	105.80
26	14	929	G	C5-C6-O6	-6.80	124.52	128.60
26	14	2249	U	N3-C2-O2	-6.80	117.44	122.20
26	1H	768	G	OP1-P-OP2	6.80	129.80	119.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	2248	C	OP1-P-O3'	6.80	120.17	105.20
1	13	365	U	C5-C6-N1	6.80	126.10	122.70
26	1H	635	C	O5'-P-OP2	-6.80	99.58	105.70
26	1H	676	A	C4-C5-N7	6.80	114.10	110.70
1	1G	1529	G	N3-C4-C5	-6.80	125.20	128.60
26	14	2581	G	C8-N9-C4	-6.80	103.68	106.40
1	1G	28	G	C8-N9-C4	-6.80	103.68	106.40
23	2L	35	C	N1-C2-O2	6.80	122.98	118.90
26	1H	2062	A	N7-C8-N9	-6.80	110.40	113.80
26	1H	2681	C	C6-N1-C2	-6.80	117.58	120.30
1	13	777	A	O5'-P-OP1	6.80	118.86	110.70
26	14	1379	A	C4-C5-N7	6.80	114.10	110.70
1	13	760	G	C5-C6-O6	-6.79	124.52	128.60
26	1H	138	G	C5-N7-C8	-6.79	100.90	104.30
26	14	2056	G	O5'-P-OP1	-6.79	99.59	105.70
1	13	1502	A	N7-C8-N9	6.79	117.20	113.80
26	1H	1312	U	O5'-P-OP2	6.79	118.85	110.70
26	14	1401	G	C8-N9-C4	-6.79	103.68	106.40
26	1H	692	C	N3-C4-N4	6.79	122.75	118.00
26	14	2338	G	O5'-P-OP1	-6.79	99.59	105.70
26	1H	1831	G	C8-N9-C4	-6.79	103.69	106.40
26	14	2401	U	C5-C6-N1	6.78	126.09	122.70
1	13	47	C	N1-C2-O2	-6.78	114.83	118.90
1	13	892	A	C2-N3-C4	-6.78	107.21	110.60
26	1H	1373	A	O5'-P-OP1	6.78	118.84	110.70
26	1H	1925	C	C6-N1-C2	-6.78	117.59	120.30
41	C8	12	ARG	NE-CZ-NH2	-6.78	116.91	120.30
26	14	1474	C	C6-N1-C2	-6.78	117.59	120.30
26	1H	528	A	O4'-C1'-N9	-6.78	102.78	108.20
26	1H	2246	G	C5-N7-C8	6.78	107.69	104.30
1	1G	690	G	N7-C8-N9	6.78	116.49	113.10
26	1H	1303	G	C5-C6-O6	6.77	132.66	128.60
26	1H	1899	G	OP2-P-O3'	6.77	120.10	105.20
26	14	2429	G	OP1-P-OP2	-6.77	109.44	119.60
26	1H	1534	G	C2-N3-C4	6.77	115.28	111.90
26	1H	1931	U	C2-N3-C4	-6.77	122.94	127.00
26	1H	2210	G	C8-N9-C4	-6.77	103.69	106.40
26	14	805	G	N3-C4-N9	6.77	130.06	126.00
26	1H	1634	A	C4-C5-C6	6.77	120.38	117.00
26	1H	2616	C	O5'-P-OP2	6.77	118.82	110.70
26	14	1660	C	OP1-P-OP2	6.77	129.75	119.60
26	1H	788	A	N1-C6-N6	6.76	122.66	118.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	1477	A	O5'-P-OP2	-6.76	99.61	105.70
1	1G	108	G	C4-C5-N7	6.76	113.51	110.80
26	14	2818	G	C8-N9-C4	6.76	109.11	106.40
26	1H	2578	G	C8-N9-C4	6.76	109.11	106.40
1	13	1113	C	C6-N1-C2	-6.76	117.60	120.30
26	1H	1817	G	C8-N9-C4	6.76	109.10	106.40
26	1H	2444	G	N9-C4-C5	6.76	108.10	105.40
26	14	1332	G	C5-C6-O6	-6.76	124.55	128.60
26	1H	131	G	C5-C6-O6	-6.76	124.55	128.60
26	1H	468	G	C5-C6-O6	-6.76	124.55	128.60
26	1H	1299	G	N1-C6-O6	6.76	123.95	119.90
26	14	693	C	C5-C6-N1	-6.76	117.62	121.00
26	1H	678	C	C5-C6-N1	-6.75	117.62	121.00
26	1H	579	G	N1-C2-N3	-6.75	119.85	123.90
1	1G	108	G	C5-C6-O6	-6.75	124.55	128.60
26	14	570	G	C4-C5-C6	6.75	122.85	118.80
26	1H	404	C	C6-N1-C2	6.75	123.00	120.30
26	1H	2379	G	C5-C6-O6	-6.75	124.55	128.60
26	14	784	A	C5-C6-N1	-6.75	114.32	117.70
26	14	1826	G	N7-C8-N9	-6.75	109.72	113.10
26	1H	1935	G	N3-C2-N2	-6.75	115.18	119.90
26	1H	2447	G	C6-N1-C2	-6.75	121.05	125.10
4	3E	12	CYS	CA-CB-SG	6.75	126.14	114.00
26	1H	514	A	C5-N7-C8	6.75	107.27	103.90
26	1H	874	G	O5'-P-OP2	-6.75	99.63	105.70
26	1H	1763	G	O5'-P-OP2	-6.74	99.63	105.70
1	13	816	A	N9-C4-C5	6.74	108.50	105.80
26	1H	121	G	N3-C4-N9	6.74	130.04	126.00
26	1H	481	G	O5'-P-OP2	-6.74	99.63	105.70
26	1H	1653	G	N3-C2-N2	6.74	124.62	119.90
26	1H	2032	G	C2-N3-C4	-6.74	108.53	111.90
26	1H	2318	G	C8-N9-C4	-6.74	103.70	106.40
26	1H	2467	C	C5-C6-N1	-6.74	117.63	121.00
26	1H	1160	G	C8-N9-C4	-6.74	103.70	106.40
26	1H	600	G	N3-C4-C5	6.74	131.97	128.60
26	1H	860	U	N1-C2-N3	6.74	118.94	114.90
26	14	451	C	OP1-P-OP2	6.74	129.71	119.60
1	13	1496	C	C6-N1-C1'	6.74	128.88	120.80
26	1H	1443	G	C4-N9-C1'	6.74	135.26	126.50
26	1H	2563	U	O5'-P-OP1	-6.74	99.64	105.70
1	13	1521	G	O5'-P-OP1	-6.73	99.64	105.70
1	1G	690	G	O4'-C1'-N9	6.73	113.59	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	13	1290	G	C8-N9-C4	-6.73	103.71	106.40
26	1H	1265	A	N7-C8-N9	6.73	117.17	113.80
26	14	935	C	C6-N1-C2	6.73	122.99	120.30
26	14	2079	U	O5'-P-OP1	-6.73	99.64	105.70
1	13	970	C	O5'-P-OP1	-6.73	99.64	105.70
26	14	2506	U	C2-N1-C1'	6.73	125.78	117.70
26	1H	1204	A	N1-C6-N6	6.73	122.64	118.60
26	1H	1307	A	N9-C4-C5	-6.73	103.11	105.80
26	1H	1614	A	C5-C6-N1	-6.73	114.33	117.70
26	1H	372	G	C8-N9-C4	-6.73	103.71	106.40
26	1H	1678	G	N3-C4-C5	6.73	131.96	128.60
25	4L	21	C	C6-N1-C2	-6.73	117.61	120.30
26	1H	245	G	N3-C4-N9	6.72	130.03	126.00
26	14	472	A	N9-C4-C5	6.72	108.49	105.80
26	1H	197	A	N1-C2-N3	6.72	132.66	129.30
26	1H	690	G	N1-C6-O6	6.72	123.93	119.90
26	1H	917	A	C5-C6-N1	-6.72	114.34	117.70
26	14	974(A)	C	N1-C2-O2	6.72	122.93	118.90
26	1H	573	G	C2-N3-C4	6.72	115.26	111.90
26	14	775	G	N1-C6-O6	-6.72	115.87	119.90
26	1H	704	G	C8-N9-C4	-6.72	103.71	106.40
26	1H	1700	A	OP1-P-OP2	6.72	129.68	119.60
26	1H	1885	A	C8-N9-C4	6.72	108.49	105.80
26	1H	1971	A	C5-C6-N1	6.72	121.06	117.70
26	1H	655	A	N1-C6-N6	6.71	122.63	118.60
26	14	1950	G	N3-C4-N9	6.71	130.03	126.00
26	1H	790	C	N1-C2-O2	-6.71	114.87	118.90
26	14	1278	A	C2-N3-C4	-6.71	107.24	110.60
26	14	1613	G	N1-C2-N2	-6.71	110.16	116.20
26	1H	1577	C	C6-N1-C2	-6.71	117.62	120.30
27	16	79	C	N3-C4-N4	6.71	122.70	118.00
26	14	783	A	N7-C8-N9	6.71	117.15	113.80
26	14	2610	C	N3-C4-C5	6.71	124.58	121.90
1	1G	230	G	C5-C6-N1	-6.71	108.15	111.50
26	1H	1982	C	C2-N3-C4	6.70	123.25	119.90
26	1H	778	G	O5'-P-OP2	-6.70	99.67	105.70
26	1H	189	G	N3-C2-N2	-6.70	115.21	119.90
26	14	1999	C	OP2-P-O3'	6.70	119.94	105.20
24	3K	76	A	C2-N3-C4	-6.70	107.25	110.60
26	14	1999	C	N3-C4-C5	6.70	124.58	121.90
26	14	2252	G	O5'-P-OP2	-6.70	99.67	105.70
26	14	1258	C	N3-C4-C5	6.69	124.58	121.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	2584	U	C2-N1-C1'	6.69	125.73	117.70
26	1H	1249	U	O5'-P-OP2	6.69	118.73	110.70
26	1H	77	C	OP1-P-OP2	-6.69	109.56	119.60
26	14	2501	C	P-O3'-C3'	6.69	127.73	119.70
26	14	464	U	O5'-P-OP2	-6.68	99.68	105.70
26	14	676	A	N1-C2-N3	6.68	132.64	129.30
1	13	776	G	O5'-P-OP1	-6.68	99.69	105.70
26	1H	210	C	C6-N1-C2	6.67	122.97	120.30
26	1H	560	C	N1-C2-O2	-6.67	114.90	118.90
26	1H	1931	U	C5-C4-O4	6.67	129.90	125.90
27	16	44	G	C4-N9-C1'	-6.67	117.83	126.50
26	14	621	A	C5-N7-C8	-6.67	100.56	103.90
26	14	2439	A	C5-N7-C8	-6.67	100.57	103.90
26	1H	67	U	C6-N1-C2	-6.67	117.00	121.00
26	1H	632	A	N7-C8-N9	6.67	117.13	113.80
26	14	48	G	OP2-P-O3'	6.67	119.86	105.20
26	14	1348	G	O5'-P-OP2	6.67	118.70	110.70
26	1H	1695	G	O5'-P-OP1	-6.66	99.70	105.70
26	1H	1828	G	C2-N3-C4	-6.66	108.57	111.90
26	1H	2053	G	N3-C2-N2	-6.66	115.24	119.90
26	14	1204	A	C2-N3-C4	-6.66	107.27	110.60
27	1J	30	C	C6-N1-C2	-6.66	117.64	120.30
26	1H	1569	A	N7-C8-N9	6.66	117.13	113.80
26	14	2326	C	O5'-P-OP1	-6.65	99.71	105.70
26	1H	2353	G	O5'-P-OP1	-6.65	99.71	105.70
26	1H	2731	G	C5-C6-O6	-6.65	124.61	128.60
26	1H	1380	G	C6-C5-N7	-6.65	126.41	130.40
26	14	1304	C	N3-C2-O2	-6.65	117.25	121.90
26	14	2314	C	C6-N1-C2	-6.65	117.64	120.30
26	14	208	C	C5-C4-N4	-6.65	115.55	120.20
26	1H	2281	C	OP1-P-O3'	6.65	119.82	105.20
26	1H	1379	A	N9-C4-C5	-6.64	103.14	105.80
1	13	792	A	N9-C1'-C2'	6.64	122.63	114.00
26	1H	1379	A	C5-C6-N6	-6.64	118.39	123.70
26	1H	122	G	O5'-P-OP2	-6.64	99.72	105.70
26	1H	1410	G	C4-N9-C1'	-6.64	117.87	126.50
26	1H	842	G	O5'-P-OP2	6.64	118.67	110.70
26	1H	1352	U	O5'-P-OP2	-6.64	99.72	105.70
26	14	1769	G	N3-C4-N9	6.64	129.98	126.00
26	1H	803	U	N1-C2-N3	6.64	118.88	114.90
26	1H	992	C	OP1-P-O3'	6.64	119.80	105.20
48	F5	82	LEU	CA-CB-CG	6.64	130.56	115.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	141	A	C8-N9-C4	-6.63	103.15	105.80
26	1H	1632	A	C5-N7-C8	-6.63	100.58	103.90
26	1H	2269	A	C8-N9-C4	6.63	108.45	105.80
26	1H	2447	G	O4'-C1'-N9	6.63	113.51	108.20
26	1H	566	U	C5-C4-O4	-6.63	121.92	125.90
26	1H	2018	G	C4-C5-N7	6.63	113.45	110.80
26	1H	2869	G	C8-N9-C4	-6.63	103.75	106.40
26	1H	871	U	N3-C4-O4	6.63	124.04	119.40
26	1H	973	A	C2-N3-C4	-6.63	107.28	110.60
26	1H	966	G	N3-C2-N2	6.63	124.54	119.90
26	14	2702	U	N3-C2-O2	-6.63	117.56	122.20
26	1H	866	A	N9-C4-C5	-6.63	103.15	105.80
26	1H	1993	U	C2-N3-C4	-6.62	123.03	127.00
26	1H	2555	U	N1-C2-N3	6.62	118.87	114.90
26	1H	1248	G	N1-C2-N2	6.62	122.16	116.20
26	14	691	C	N3-C4-C5	-6.62	119.25	121.90
1	13	311	C	C5-C6-N1	6.62	124.31	121.00
26	14	492	A	O5'-P-OP2	-6.62	99.74	105.70
26	1H	778	G	N1-C6-O6	-6.62	115.93	119.90
26	1H	2518	A	C5-C6-N6	-6.62	118.41	123.70
26	14	1251	C	N3-C4-N4	6.62	122.63	118.00
26	1H	177	G	N1-C6-O6	-6.61	115.93	119.90
26	1H	1425	G	C2-N3-C4	6.61	115.21	111.90
26	1H	1613	G	N3-C2-N2	6.61	124.53	119.90
26	14	140	A	C8-N9-C4	-6.61	103.16	105.80
26	14	1371	G	N1-C6-O6	6.61	123.87	119.90
26	1H	265	A	C5-C6-N1	-6.61	114.39	117.70
26	1H	2390	U	O5'-P-OP1	-6.61	99.75	105.70
26	14	74	A	C4-C5-N7	6.61	114.01	110.70
26	1H	1248	G	N3-C2-N2	-6.61	115.27	119.90
1	13	690	G	N7-C8-N9	6.61	116.41	113.10
26	1H	2713	A	C4-C5-N7	6.61	114.00	110.70
1	1G	972	C	C6-N1-C2	-6.61	117.66	120.30
26	1H	1496	A	C6-C5-N7	-6.61	127.68	132.30
26	1H	1204	A	C4-N9-C1'	6.60	138.19	126.30
1	1G	197	A	P-O3'-C3'	6.60	127.62	119.70
1	13	812	C	P-O3'-C3'	6.60	127.62	119.70
26	1H	944	G	C5-C6-O6	6.60	132.56	128.60
26	1H	1332	G	C5-C6-N1	-6.60	108.20	111.50
26	1H	2446	G	N9-C4-C5	-6.60	102.76	105.40
26	14	1342	A	C6-C5-N7	-6.60	127.68	132.30
26	1H	216	A	O5'-P-OP2	6.60	118.62	110.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	872	A	O5'-P-OP1	-6.60	99.76	105.70
26	14	786	C	O5'-P-OP2	-6.60	99.76	105.70
26	14	2712	U	C5-C4-O4	6.60	129.86	125.90
26	1H	1776	G	O5'-P-OP1	6.60	118.62	110.70
26	1H	1806	C	O5'-P-OP2	-6.60	99.76	105.70
28	11	95	LEU	CA-CB-CG	6.60	130.47	115.30
24	1L	74	C	C6-N1-C2	6.60	122.94	120.30
1	1G	1339	A	O4'-C1'-N9	6.60	113.48	108.20
1	13	715	A	O5'-P-OP2	-6.59	99.77	105.70
1	1G	114	U	C5-C6-N1	-6.59	119.40	122.70
24	1L	74	C	O4'-C1'-N1	6.59	113.47	108.20
26	1H	122	G	C2-N3-C4	-6.59	108.61	111.90
26	1H	1681	G	N3-C4-C5	6.59	131.90	128.60
26	14	1686	C	C6-N1-C2	6.59	122.94	120.30
27	1J	70	C	C6-N1-C2	-6.59	117.67	120.30
26	1H	509	C	OP2-P-O3'	6.59	119.69	105.20
40	B8	96	ARG	NE-CZ-NH1	-6.59	117.01	120.30
1	1G	912	C	O5'-P-OP2	-6.59	99.77	105.70
26	14	210	C	N3-C4-N4	-6.59	113.39	118.00
1	13	1219	U	C5-C6-N1	6.59	125.99	122.70
26	1H	1977	A	C8-N9-C4	6.59	108.43	105.80
26	1H	2301	C	C6-N1-C2	-6.59	117.67	120.30
26	1H	2385	C	O5'-P-OP2	-6.59	99.77	105.70
1	13	23	C	C5-C6-N1	6.58	124.29	121.00
26	14	1277	G	C8-N9-C4	6.58	109.03	106.40
1	1G	1529	G	C4-N9-C1'	6.58	135.06	126.50
26	14	2011	U	O5'-P-OP1	-6.58	99.78	105.70
26	1H	141	A	O4'-C1'-N9	6.58	113.47	108.20
26	1H	779	U	N1-C2-N3	-6.58	110.95	114.90
26	1H	808	G	N1-C2-N3	6.58	127.85	123.90
26	1H	2096	U	C6-N1-C2	-6.58	117.05	121.00
26	14	212	G	O5'-P-OP2	-6.58	99.78	105.70
26	14	395	U	O4'-C1'-N1	6.58	113.46	108.20
26	1H	582	G	C6-C5-N7	-6.58	126.45	130.40
26	1H	2703	C	N3-C2-O2	-6.58	117.30	121.90
26	14	2013	A	C2-N3-C4	-6.58	107.31	110.60
26	1H	58	G	C8-N9-C4	-6.58	103.77	106.40
26	1H	945	A	C4-C5-N7	6.58	113.99	110.70
26	1H	2013	A	N1-C6-N6	-6.58	114.66	118.60
1	1G	1286	A	N7-C8-N9	6.58	117.09	113.80
26	14	1621	U	N1-C2-O2	-6.58	118.20	122.80
26	14	2506	U	O4'-C1'-N1	-6.58	102.94	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	128	C	C6-N1-C2	6.57	122.93	120.30
26	1H	355	G	N1-C6-O6	6.57	123.84	119.90
26	14	2056	G	N1-C6-O6	6.57	123.84	119.90
26	1H	1978	A	O5'-P-OP1	-6.57	99.79	105.70
1	1G	11	G	O5'-P-OP1	-6.57	99.79	105.70
26	14	1791	A	OP1-P-OP2	-6.57	109.75	119.60
26	14	2256	G	N3-C2-N2	6.57	124.50	119.90
26	1H	2324	C	N3-C4-C5	6.57	124.53	121.90
26	1H	577	G	OP1-P-OP2	-6.56	109.75	119.60
26	1H	1786	A	C4-N9-C1'	6.56	138.11	126.30
26	14	93	C	C5-C6-N1	6.56	124.28	121.00
26	14	2840	C	O5'-P-OP2	-6.56	99.79	105.70
1	13	623	C	C6-N1-C2	-6.56	117.68	120.30
26	1H	599	G	N3-C2-N2	6.56	124.49	119.90
26	14	929	G	C8-N9-C4	-6.56	103.78	106.40
26	1H	1348	G	OP1-P-O3'	6.56	119.62	105.20
26	14	148	C	C6-N1-C2	6.56	122.92	120.30
26	1H	774	A	C4-N9-C1'	-6.55	114.50	126.30
26	1H	1931	U	C4-C5-C6	6.55	123.63	119.70
26	1H	1781	C	C2-N1-C1'	6.55	126.01	118.80
26	1H	148	C	C6-N1-C2	6.55	122.92	120.30
26	1H	381	G	N1-C6-O6	-6.55	115.97	119.90
26	1H	787	U	O5'-P-OP1	6.55	118.56	110.70
26	1H	954	G	O5'-P-OP2	6.55	118.56	110.70
26	14	828	U	N3-C2-O2	-6.55	117.61	122.20
26	14	1325	G	N9-C4-C5	-6.55	102.78	105.40
1	13	977	A	N1-C6-N6	-6.55	114.67	118.60
1	13	1508	G	N3-C4-N9	-6.55	122.07	126.00
26	1H	1623	G	OP2-P-O3'	6.55	119.60	105.20
26	1H	2871	C	O5'-P-OP1	6.55	118.56	110.70
26	1H	127	A	N1-C6-N6	6.54	122.53	118.60
26	1H	1618	A	OP1-P-OP2	-6.54	109.78	119.60
26	1H	2228	G	C8-N9-C1'	-6.54	118.49	127.00
26	14	1187	G	C8-N9-C4	-6.54	103.78	106.40
26	1H	482	A	C8-N9-C4	-6.54	103.18	105.80
26	1H	763	G	OP2-P-O3'	6.54	119.59	105.20
26	1H	2074	U	O5'-P-OP1	-6.54	99.81	105.70
1	1G	1498	U	C2-N1-C1'	6.54	125.55	117.70
26	1H	910	A	N1-C6-N6	6.54	122.52	118.60
26	14	2401	U	C2-N1-C1'	6.54	125.55	117.70
26	1H	454	A	O5'-P-OP2	-6.54	99.82	105.70
26	1H	663	G	N3-C4-C5	-6.54	125.33	128.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	1375	C	OP1-P-O3'	6.54	119.58	105.20
26	1H	1616	A	OP1-P-OP2	6.54	129.41	119.60
26	14	1823	G	C5-C6-O6	6.54	132.52	128.60
26	14	2426	A	C5-N7-C8	-6.54	100.63	103.90
26	1H	528	A	C8-N9-C1'	6.53	139.46	127.70
26	1H	663	G	OP1-P-OP2	6.53	129.40	119.60
26	1H	2271	G	N3-C4-C5	-6.53	125.33	128.60
1	13	963	G	N1-C2-N2	-6.53	110.32	116.20
1	13	1205	U	N3-C4-C5	-6.53	110.68	114.60
26	1H	1157	G	C2-N3-C4	-6.53	108.64	111.90
26	1H	1265	A	N1-C2-N3	6.53	132.56	129.30
26	1H	1667	G	N3-C4-C5	6.53	131.87	128.60
1	1G	330	C	N1-C2-O2	6.53	122.82	118.90
26	1H	1966	A	O5'-P-OP2	-6.53	99.83	105.70
26	1H	1157	G	N1-C2-N3	6.53	127.82	123.90
26	1H	2331	G	N9-C4-C5	-6.53	102.79	105.40
26	14	1840	G	N3-C2-N2	-6.53	115.33	119.90
26	14	2244	U	N3-C2-O2	6.53	126.77	122.20
1	13	1498	U	P-O3'-C3'	6.52	127.53	119.70
26	1H	1574	C	C5-C6-N1	-6.52	117.74	121.00
26	14	2740	A	O5'-P-OP2	-6.52	99.83	105.70
26	1H	22	C	O5'-P-OP1	-6.52	99.83	105.70
26	1H	1128	A	OP1-P-OP2	6.52	129.38	119.60
26	1H	1559	G	N3-C4-C5	6.52	131.86	128.60
1	13	578	C	N3-C4-C5	-6.52	119.29	121.90
26	1H	688	U	O5'-P-OP2	-6.52	99.84	105.70
26	1H	2346	A	C4-N9-C1'	6.51	138.03	126.30
26	14	832	G	C8-N9-C4	-6.51	103.79	106.40
26	1H	1204	A	C5-C6-N1	-6.51	114.44	117.70
26	1H	821	A	O5'-P-OP2	-6.51	99.84	105.70
26	14	2243	U	OP2-P-O3'	6.51	119.52	105.20
26	14	2688	U	C5-C4-O4	6.51	129.81	125.90
26	1H	1623	G	N1-C6-O6	-6.51	116.00	119.90
26	1H	1674	G	C4-N9-C1'	6.50	134.96	126.50
1	13	1266	G	C6-C5-N7	6.50	134.30	130.40
26	1H	1278	A	C8-N9-C4	6.50	108.40	105.80
26	1H	2228	G	N3-C4-N9	6.50	129.90	126.00
26	1H	51	G	C5-C6-O6	6.50	132.50	128.60
26	1H	1970	A	C8-N9-C4	-6.50	103.20	105.80
26	1H	2342	C	C6-N1-C2	-6.50	117.70	120.30
26	14	805	G	C5-C6-O6	-6.50	124.70	128.60
26	14	76	C	C6-N1-C2	-6.50	117.70	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	1969	A	C5-N7-C8	6.49	107.15	103.90
1	1G	353	A	N1-C6-N6	6.49	122.50	118.60
26	14	388	G	N3-C4-C5	6.49	131.85	128.60
26	1H	834	C	C6-N1-C2	-6.49	117.70	120.30
26	14	2713	A	C8-N9-C4	-6.49	103.20	105.80
26	1H	1397	U	N3-C2-O2	-6.49	117.66	122.20
26	1H	2263	C	N3-C4-C5	-6.49	119.30	121.90
26	14	455	C	N1-C2-O2	6.49	122.79	118.90
26	1H	1417	C	N3-C4-C5	-6.49	119.31	121.90
26	1H	579	G	C2-N3-C4	6.49	115.14	111.90
26	1H	1308	A	C8-N9-C4	-6.49	103.21	105.80
30	31	176	LEU	CB-CG-CD2	-6.49	99.98	111.00
26	1H	2702	U	C6-N1-C2	-6.48	117.11	121.00
26	1H	1831	G	OP1-P-OP2	-6.48	109.88	119.60
26	14	566	U	C6-N1-C2	6.48	124.89	121.00
26	1H	698	C	C6-N1-C2	6.48	122.89	120.30
26	14	138	G	O4'-C1'-N9	6.48	113.38	108.20
26	1H	210	C	N3-C4-C5	6.48	124.49	121.90
26	1H	467	G	N7-C8-N9	-6.48	109.86	113.10
26	14	1570	A	N1-C6-N6	6.48	122.49	118.60
26	14	2263	C	C5-C6-N1	6.48	124.24	121.00
26	1H	1621	U	C4-C5-C6	6.48	123.58	119.70
26	14	511	U	C6-N1-C2	-6.47	117.11	121.00
1	13	664	G	O5'-P-OP2	-6.47	99.87	105.70
26	1H	2330	G	C2-N3-C4	-6.47	108.66	111.90
27	16	8	U	O5'-P-OP1	6.47	118.47	110.70
26	1H	74	A	N3-C4-N9	-6.47	122.22	127.40
26	1H	952	G	O5'-P-OP2	6.47	118.46	110.70
26	1H	1425	G	C8-N9-C4	-6.47	103.81	106.40
26	1H	2228	G	N1-C6-O6	6.47	123.78	119.90
26	14	121	G	N1-C6-O6	6.47	123.78	119.90
26	14	270(Q)	C	C6-N1-C2	-6.47	117.71	120.30
26	1H	247	G	C8-N9-C4	6.47	108.99	106.40
26	1H	1248	G	N1-C6-O6	6.47	123.78	119.90
26	1H	2286	A	N7-C8-N9	6.47	117.03	113.80
27	16	9	G	OP2-P-O3'	6.47	119.43	105.20
26	14	2075	U	C5-C6-N1	-6.47	119.47	122.70
26	1H	85	G	O5'-P-OP2	-6.46	99.88	105.70
26	1H	1443	G	N1-C6-O6	6.46	123.78	119.90
26	1H	2713	A	N1-C6-N6	6.46	122.48	118.60
27	16	12	C	C5-C6-N1	-6.46	117.77	121.00
24	3K	71	G	C4-N9-C1'	-6.46	118.10	126.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	2449	U	N1-C2-N3	6.46	118.78	114.90
26	14	775	G	N3-C4-N9	6.46	129.88	126.00
26	1H	613	U	N3-C4-O4	-6.46	114.88	119.40
26	1H	1786	A	N9-C1'-C2'	6.46	122.40	114.00
1	1G	1192	C	C6-N1-C2	-6.46	117.72	120.30
26	1H	192	C	N3-C2-O2	6.46	126.42	121.90
26	14	2297	C	O5'-P-OP1	-6.46	99.89	105.70
26	14	2427	C	C5-C4-N4	-6.46	115.68	120.20
26	1H	1660	C	N3-C4-N4	-6.46	113.48	118.00
1	1G	519	C	C6-N1-C2	6.46	122.88	120.30
32	51	172	LYS	C-N-CD	-6.46	106.40	120.60
26	1H	632	A	C8-N9-C4	-6.45	103.22	105.80
9	8E	47	LEU	CA-CB-CG	6.45	130.14	115.30
26	14	579	G	N1-C6-O6	6.45	123.77	119.90
1	13	1266	G	C8-N9-C1'	6.45	135.38	127.00
26	14	959	A	OP1-P-O3'	6.45	119.39	105.20
26	1H	1781	C	N3-C4-C5	6.45	124.48	121.90
26	14	494	G	N1-C6-O6	6.45	123.77	119.90
26	14	580	C	N3-C4-C5	-6.45	119.32	121.90
26	14	1574	C	C5-C4-N4	-6.45	115.69	120.20
26	14	1608	A	C5-C6-N6	6.45	128.86	123.70
26	14	1808	U	O5'-P-OP1	-6.45	99.90	105.70
26	1H	116	C	N3-C4-C5	-6.44	119.32	121.90
26	14	1999	C	C6-N1-C2	6.44	122.88	120.30
1	13	186	C	C6-N1-C2	-6.44	117.72	120.30
26	1H	2263	C	C5-C4-N4	6.44	124.71	120.20
26	1H	411	G	N9-C4-C5	6.44	107.98	105.40
26	1H	1938	A	C4-C5-C6	6.44	120.22	117.00
26	14	2042	A	O5'-P-OP2	-6.44	99.90	105.70
26	14	2071	A	C6-N1-C2	-6.44	114.73	118.60
26	14	2505	G	C5-C6-N1	-6.44	108.28	111.50
26	1H	271(B)	G	N3-C4-N9	6.44	129.86	126.00
26	14	1313	U	C5-C6-N1	6.44	125.92	122.70
26	1H	955	C	N3-C4-C5	-6.43	119.33	121.90
26	14	593	G	C8-N9-C4	6.43	108.97	106.40
26	1H	114	U	OP1-P-O3'	6.43	119.35	105.20
26	1H	2313	C	OP2-P-O3'	6.43	119.35	105.20
27	16	109	G	C8-N9-C4	-6.43	103.83	106.40
1	1G	818	G	C5-C6-O6	6.43	132.46	128.60
26	1H	190	A	C5-C6-N1	6.43	120.92	117.70
26	1H	1670	C	C5-C6-N1	-6.43	117.78	121.00
26	14	2263	C	C6-N1-C2	-6.43	117.73	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	1558	A	N1-C2-N3	6.43	132.51	129.30
37	88	86	GLY	N-CA-C	-6.43	97.03	113.10
1	13	1051	C	C6-N1-C2	-6.42	117.73	120.30
26	1H	2374	C	O5'-P-OP2	-6.42	99.92	105.70
26	1H	2380	C	O5'-P-OP2	-6.42	99.92	105.70
26	1H	2726	U	N3-C4-O4	-6.42	114.90	119.40
26	14	1815	A	OP1-P-O3'	6.42	119.33	105.20
26	14	1950	G	C2-N3-C4	6.42	115.11	111.90
26	1H	265	A	O4'-C1'-N9	6.42	113.34	108.20
26	14	1332	G	N1-C2-N2	-6.42	110.42	116.20
26	14	1408	C	N1-C2-O2	-6.42	115.05	118.90
26	1H	213	A	C4-C5-C6	-6.42	113.79	117.00
26	1H	866	A	N1-C6-N6	6.42	122.45	118.60
26	1H	1590	U	O5'-P-OP1	-6.42	99.92	105.70
26	1H	1398	C	O5'-P-OP2	6.42	118.40	110.70
26	1H	746	A	O5'-P-OP1	-6.42	99.92	105.70
26	14	1613	G	C8-N9-C4	6.42	108.97	106.40
26	14	916	G	C8-N9-C4	-6.42	103.83	106.40
26	14	1613	G	C8-N9-C1'	-6.42	118.66	127.00
27	16	115	G	C6-N1-C2	-6.42	121.25	125.10
1	1G	723	U	C2-N1-C1'	6.42	125.40	117.70
26	1H	609	A	C8-N9-C4	6.41	108.37	105.80
26	1H	1596	A	OP2-P-O3'	6.41	119.31	105.20
1	1G	354	G	C4-N9-C1'	6.41	134.84	126.50
26	14	1607	C	N3-C4-N4	6.41	122.49	118.00
26	1H	182	A	C8-N9-C4	6.41	108.36	105.80
24	3L	76	A	C4-C5-N7	6.41	113.91	110.70
1	13	22	G	N3-C2-N2	-6.41	115.42	119.90
26	1H	594	U	C6-N1-C2	6.41	124.84	121.00
26	14	2039	C	N1-C2-O2	6.41	122.75	118.90
26	1H	672	C	O5'-P-OP2	-6.41	99.93	105.70
26	1H	444	C	OP1-P-OP2	-6.41	109.99	119.60
26	1H	1373	A	N7-C8-N9	-6.41	110.60	113.80
26	1H	1600	C	OP1-P-O3'	6.41	119.29	105.20
1	13	768	A	C8-N9-C4	-6.40	103.24	105.80
26	1H	2072	G	N9-C4-C5	-6.40	102.84	105.40
26	1H	2494	G	C5-C6-N1	-6.40	108.30	111.50
26	1H	2713	A	C5-C6-N1	-6.40	114.50	117.70
26	14	1784	A	C5-N7-C8	-6.40	100.70	103.90
1	13	910	C	C6-N1-C2	6.40	122.86	120.30
1	13	963	G	N3-C4-N9	6.40	129.84	126.00
26	14	1375	C	OP1-P-O3'	6.40	119.28	105.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	13	899	C	N3-C2-O2	6.40	126.38	121.90
26	14	775	G	C5-C6-N1	6.40	114.70	111.50
26	14	1318	C	C6-N1-C2	-6.40	117.74	120.30
1	13	1279	A	N7-C8-N9	6.39	117.00	113.80
26	1H	24	G	C5-C6-N1	-6.39	108.30	111.50
26	1H	2241	A	C2-N3-C4	-6.39	107.40	110.60
26	14	1594	G	C8-N9-C4	-6.39	103.84	106.40
26	14	2598	A	C8-N9-C4	6.39	108.36	105.80
26	1H	2552	U	N1-C2-N3	6.39	118.73	114.90
26	14	194	G	O5'-P-OP2	6.39	118.37	110.70
26	14	774	A	C4-N9-C1'	-6.39	114.79	126.30
26	1H	468	G	OP1-P-OP2	-6.39	110.02	119.60
26	14	2439	A	C8-N9-C4	-6.39	103.24	105.80
26	1H	1204	A	N1-C2-N3	6.39	132.49	129.30
26	14	1995	U	O5'-P-OP2	-6.39	99.95	105.70
23	2L	21	U	N3-C2-O2	-6.39	117.73	122.20
26	14	372	G	O4'-C1'-N9	6.39	113.31	108.20
26	14	821	A	C5-C6-N6	-6.39	118.59	123.70
26	14	2451	A	C8-N9-C4	-6.39	103.25	105.80
26	1H	2395	C	C5-C4-N4	-6.38	115.73	120.20
26	1H	2467	C	C6-N1-C2	6.38	122.85	120.30
26	1H	663	G	O5'-P-OP2	-6.38	99.95	105.70
26	1H	1355	G	N1-C6-O6	-6.38	116.07	119.90
26	1H	2412	A	C6-N1-C2	-6.38	114.77	118.60
26	1H	94	G	C5-C6-O6	-6.38	124.77	128.60
26	1H	1575	C	O5'-P-OP1	6.38	118.36	110.70
26	1H	2379	G	C8-N9-C1'	-6.38	118.70	127.00
26	1H	2358	G	C8-N9-C4	6.38	108.95	106.40
26	1H	685	A	C5-N7-C8	-6.38	100.71	103.90
26	1H	1280	G	OP1-P-OP2	-6.38	110.03	119.60
26	1H	2025	C	C6-N1-C2	-6.38	117.75	120.30
1	1G	1417	G	C6-C5-N7	-6.38	126.57	130.40
26	14	856	C	C5-C6-N1	6.38	124.19	121.00
26	14	1135	C	N1-C2-O2	6.38	122.73	118.90
26	14	2281	C	O5'-P-OP2	6.38	118.35	110.70
26	14	2572	A	O5'-P-OP1	-6.38	99.96	105.70
1	13	966	G	N9-C4-C5	-6.38	102.85	105.40
26	1H	958	U	O5'-P-OP1	6.38	118.35	110.70
1	1G	1346	A	OP2-P-O3'	6.38	119.23	105.20
26	14	1443	G	N1-C6-O6	6.38	123.72	119.90
26	14	2681	C	N3-C4-N4	-6.38	113.54	118.00
26	1H	120	U	N3-C2-O2	-6.37	117.74	122.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	2029	G	O5'-P-OP1	-6.37	99.96	105.70
26	1H	2618	G	N7-C8-N9	6.37	116.29	113.10
26	14	2261	C	O5'-P-OP2	-6.37	99.96	105.70
26	14	2570	G	N1-C6-O6	6.37	123.72	119.90
26	14	2713	A	N1-C6-N6	6.37	122.42	118.60
26	14	1332	G	N9-C4-C5	-6.37	102.85	105.40
26	1H	1938	A	N1-C6-N6	6.37	122.42	118.60
26	1H	1908	C	C6-N1-C2	-6.37	117.75	120.30
26	14	2615	U	O5'-P-OP2	-6.37	99.97	105.70
26	14	922	U	C6-N1-C2	-6.37	117.18	121.00
1	13	720	C	C6-N1-C2	-6.36	117.75	120.30
26	1H	2286	A	C8-N9-C4	-6.36	103.25	105.80
26	14	1154	G	N1-C6-O6	6.36	123.72	119.90
26	14	2019	A	C8-N9-C4	6.36	108.34	105.80
26	1H	853	G	O5'-P-OP2	-6.36	99.98	105.70
26	1H	2216	G	C8-N9-C4	-6.36	103.86	106.40
26	14	2610	C	O5'-P-OP1	-6.36	99.98	105.70
26	14	2763	G	N3-C4-C5	-6.36	125.42	128.60
26	1H	860	U	C2-N3-C4	-6.36	123.19	127.00
26	1H	2427	C	O5'-P-OP2	6.36	118.33	110.70
26	14	1616	A	C4-C5-N7	6.36	113.88	110.70
26	14	1824	G	N1-C6-O6	6.36	123.72	119.90
26	14	1899	G	C2-N3-C4	-6.36	108.72	111.90
26	14	470	A	OP1-P-OP2	6.36	129.13	119.60
26	1H	196	A	C5-C6-N1	-6.35	114.52	117.70
26	1H	509	C	O5'-P-OP2	-6.35	99.98	105.70
26	1H	1265	A	C4-C5-C6	6.35	120.18	117.00
26	1H	1691	C	N3-C4-C5	-6.35	119.36	121.90
26	1H	966	G	N1-C2-N2	-6.35	110.49	116.20
26	1H	2779	U	N3-C2-O2	-6.35	117.75	122.20
26	14	1527	G	N3-C2-N2	-6.35	115.45	119.90
26	1H	134	C	C5-C6-N1	-6.35	117.83	121.00
26	1H	463	G	N1-C2-N2	-6.35	110.49	116.20
26	14	1950	G	C8-N9-C4	-6.35	103.86	106.40
26	1H	764	A	O5'-P-OP1	6.34	118.31	110.70
26	1H	1936	A	C5-N7-C8	-6.34	100.73	103.90
26	1H	2308	G	C6-N1-C2	6.34	128.91	125.10
26	1H	2597	G	C5-N7-C8	-6.34	101.13	104.30
26	14	386	G	C5-C6-O6	-6.34	124.79	128.60
26	1H	391	G	C5-C6-O6	-6.34	124.79	128.60
1	1G	1527	C	C5-C6-N1	-6.34	117.83	121.00
26	14	808	G	O5'-P-OP2	-6.34	99.99	105.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	140	A	OP2-P-O3'	6.34	119.15	105.20
26	1H	692	C	C2-N3-C4	-6.34	116.73	119.90
26	1H	968	G	O5'-P-OP1	6.34	118.31	110.70
26	1H	1443	G	N1-C2-N3	6.34	127.70	123.90
26	1H	1518	C	O5'-P-OP2	6.34	118.31	110.70
26	1H	1528	A	C5-N7-C8	-6.34	100.73	103.90
27	16	99	A	O5'-P-OP2	-6.34	99.99	105.70
26	1H	1209	G	C5-C6-O6	-6.34	124.80	128.60
26	1H	1254	A	N9-C4-C5	-6.34	103.26	105.80
26	1H	1300	U	C6-N1-C1'	6.34	130.08	121.20
26	1H	38	A	OP1-P-O3'	6.34	119.14	105.20
26	1H	664	C	C5-C6-N1	-6.34	117.83	121.00
26	14	465	G	N3-C4-C5	-6.34	125.43	128.60
26	1H	769	G	C5-C6-O6	6.34	132.40	128.60
1	13	789	U	N3-C4-C5	-6.33	110.80	114.60
26	1H	679	C	C5-C6-N1	-6.33	117.83	121.00
26	1H	2513	G	N3-C4-C5	-6.33	125.43	128.60
1	1G	1417	G	N1-C6-O6	6.33	123.70	119.90
26	1H	1307	A	C8-N9-C4	6.33	108.33	105.80
26	1H	747	U	OP1-P-OP2	6.33	129.10	119.60
26	1H	1272	A	C4-C5-C6	-6.33	113.83	117.00
26	1H	690	G	N7-C8-N9	-6.33	109.94	113.10
1	13	872	A	C6-N1-C2	6.33	122.40	118.60
26	1H	138	G	O4'-C1'-N9	6.33	113.26	108.20
26	1H	2375	G	N9-C1'-C2'	-6.33	105.04	112.00
26	14	613	U	N3-C2-O2	-6.33	117.77	122.20
26	1H	1303	G	OP2-P-O3'	6.33	119.12	105.20
26	1H	1973	G	C8-N9-C4	-6.33	103.87	106.40
26	1H	797	C	C5-C6-N1	-6.33	117.84	121.00
1	13	520	A	N1-C6-N6	6.32	122.39	118.60
26	1H	828	U	C2-N1-C1'	6.32	125.29	117.70
26	1H	845	G	C2-N3-C4	-6.32	108.74	111.90
26	14	1898	U	C5-C4-O4	6.32	129.69	125.90
26	14	602	G	C4-N9-C1'	6.32	134.72	126.50
1	13	690	G	O5'-P-OP1	-6.32	100.01	105.70
26	1H	199	A	N1-C2-N3	-6.32	126.14	129.30
26	1H	199	A	O5'-P-OP1	-6.32	100.01	105.70
26	1H	148	C	C5-C6-N1	-6.32	117.84	121.00
26	1H	802	A	O5'-P-OP1	6.32	118.28	110.70
26	1H	2430	A	C6-C5-N7	-6.32	127.88	132.30
26	1H	596	G	N1-C2-N2	6.32	121.89	116.20
26	1H	860	U	O5'-P-OP1	6.32	118.28	110.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	380	U	OP1-P-OP2	6.32	129.07	119.60
1	13	584	G	N1-C2-N2	-6.32	110.52	116.20
26	1H	195	A	C5-C6-N6	-6.32	118.65	123.70
26	1H	803	U	N3-C2-O2	-6.32	117.78	122.20
26	14	2356	C	N1-C2-O2	-6.32	115.11	118.90
1	13	812	C	C2-N1-C1'	6.31	125.75	118.80
23	2K	21	U	C2-N1-C1'	6.31	125.28	117.70
26	1H	1313	U	C5-C6-N1	6.31	125.86	122.70
1	13	1374	A	C2-N3-C4	-6.31	107.44	110.60
26	1H	815	C	O5'-P-OP1	6.31	118.27	110.70
26	1H	1124	C	C4-C5-C6	6.31	120.56	117.40
26	1H	2440	C	C2-N3-C4	6.31	123.06	119.90
1	1G	1498	U	C6-N1-C2	-6.31	117.21	121.00
26	1H	1520	U	C5-C4-O4	6.31	129.69	125.90
26	14	2062	A	N9-C4-C5	-6.31	103.28	105.80
26	14	1899	G	C8-N9-C4	-6.31	103.88	106.40
26	14	2248	C	N3-C4-C5	-6.31	119.38	121.90
26	1H	1369	G	N3-C4-N9	6.31	129.78	126.00
26	1H	2319	G	N3-C4-C5	-6.31	125.45	128.60
26	1H	504	U	C2-N1-C1'	6.30	125.26	117.70
26	14	1661	G	C5-C6-O6	-6.30	124.82	128.60
26	14	1678	G	C8-N9-C4	-6.30	103.88	106.40
26	14	2374	C	N3-C4-C5	6.30	124.42	121.90
26	1H	692	C	N1-C2-O2	-6.30	115.12	118.90
26	1H	842	G	O5'-P-OP1	-6.30	100.03	105.70
26	1H	2328	A	N1-C2-N3	6.30	132.45	129.30
26	14	453	C	C6-N1-C2	6.30	122.82	120.30
26	14	579	G	N3-C2-N2	-6.30	115.49	119.90
26	1H	140	A	O4'-C1'-N9	6.30	113.24	108.20
26	1H	997	G	C8-N9-C4	6.30	108.92	106.40
26	1H	2415	G	C6-C5-N7	-6.30	126.62	130.40
26	1H	2628	C	C6-N1-C2	6.30	122.82	120.30
1	13	1322	C	C2-N1-C1'	6.30	125.73	118.80
26	1H	2083	G	N1-C6-O6	6.30	123.68	119.90
26	1H	2287	A	N1-C2-N3	6.30	132.45	129.30
26	1H	2782	G	C6-C5-N7	-6.30	126.62	130.40
1	13	176	C	C6-N1-C2	-6.29	117.78	120.30
26	1H	1326	U	OP2-P-O3'	6.29	119.05	105.20
1	13	757	U	N3-C2-O2	-6.29	117.80	122.20
26	14	201	C	O5'-P-OP2	-6.29	100.04	105.70
26	1H	202	U	N1-C2-O2	6.29	127.20	122.80
26	1H	663	G	C6-C5-N7	-6.29	126.63	130.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	1641	A	O5'-P-OP2	6.29	118.25	110.70
26	1H	2440	C	N1-C2-O2	6.29	122.67	118.90
1	1G	529	G	N1-C6-O6	6.29	123.67	119.90
26	1H	1424	G	O5'-P-OP2	-6.29	100.04	105.70
26	1H	1428	C	O5'-P-OP2	6.29	118.24	110.70
26	1H	1634	A	OP1-P-OP2	6.29	129.03	119.60
26	14	733	G	C8-N9-C4	-6.29	103.89	106.40
23	2K	35	C	C6-N1-C1'	-6.28	113.26	120.80
27	16	14	U	OP1-P-OP2	6.28	129.03	119.60
39	A8	101	LEU	CA-CB-CG	6.28	129.75	115.30
26	14	242	G	C8-N9-C4	6.28	108.91	106.40
26	14	2049	G	O5'-P-OP1	-6.28	100.05	105.70
26	14	2575	C	C2-N1-C1'	-6.28	111.89	118.80
1	13	103	C	C6-N1-C2	-6.28	117.79	120.30
1	13	386	C	C2-N1-C1'	-6.28	111.89	118.80
26	1H	852	G	OP2-P-O3'	6.28	119.02	105.20
26	14	2006	C	C5-C4-N4	-6.28	115.80	120.20
1	13	122	G	C5-C6-O6	-6.28	124.83	128.60
1	13	300	A	O5'-P-OP1	-6.28	100.05	105.70
27	16	77	U	N3-C4-C5	6.28	118.37	114.60
1	13	1199	U	N3-C2-O2	-6.28	117.81	122.20
1	1G	813	U	O5'-P-OP1	-6.28	100.05	105.70
1	1G	1487	G	N1-C6-O6	6.27	123.66	119.90
1	13	1498	U	O4'-C1'-N1	-6.27	103.18	108.20
26	1H	974(A)	C	C6-N1-C2	-6.27	117.79	120.30
26	1H	1950	G	C8-N9-C4	-6.27	103.89	106.40
26	14	470	A	C5-C6-N6	-6.27	118.68	123.70
26	14	2702	U	O5'-P-OP1	6.27	118.23	110.70
26	1H	797	C	N3-C2-O2	6.27	126.29	121.90
26	1H	2499	C	N3-C4-N4	6.27	122.39	118.00
26	14	1332	G	C5-C6-N1	-6.27	108.36	111.50
26	14	2577	A	N1-C6-N6	6.27	122.36	118.60
1	13	575	G	O4'-C1'-N9	-6.27	103.19	108.20
26	1H	640	C	OP1-P-O3'	6.27	118.99	105.20
26	1H	2228	G	C4-C5-C6	6.26	122.56	118.80
26	1H	2367	G	N7-C8-N9	6.26	116.23	113.10
26	1H	2375	G	C4-C5-N7	6.26	113.31	110.80
1	13	541	G	N1-C6-O6	6.26	123.66	119.90
26	1H	1678	G	N1-C6-O6	6.26	123.66	119.90
1	1G	1502	A	C5-N7-C8	-6.26	100.77	103.90
26	1H	585	G	N1-C6-O6	6.26	123.66	119.90
1	1G	932	C	N1-C2-O2	6.26	122.66	118.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	200	U	O5'-P-OP2	-6.26	100.07	105.70
26	14	1367	A	C2-N3-C4	-6.26	107.47	110.60
26	1H	1123	C	C5-C6-N1	-6.26	117.87	121.00
26	14	2251	G	C5-N7-C8	6.26	107.43	104.30
1	13	858	G	N1-C6-O6	-6.26	116.15	119.90
26	14	2291	U	C5-C4-O4	6.26	129.65	125.90
26	1H	1349	A	O5'-P-OP2	6.25	118.21	110.70
26	1H	2476	A	C8-N9-C4	-6.25	103.30	105.80
1	1G	267	C	N1-C2-O2	6.25	122.65	118.90
26	1H	828	U	C5-C4-O4	6.25	129.65	125.90
26	1H	1425	G	C5-C6-N1	6.25	114.63	111.50
26	1H	2324	C	C5-C4-N4	-6.25	115.82	120.20
26	14	2592	G	O5'-P-OP1	6.25	118.20	110.70
26	1H	1142(A)	A	N3-C4-N9	-6.25	122.40	127.40
1	1G	1502	A	C6-C5-N7	-6.25	127.92	132.30
1	13	1204	A	N1-C6-N6	6.25	122.35	118.60
26	14	2713	A	C4-C5-N7	6.25	113.83	110.70
26	1H	364	C	C6-N1-C2	-6.25	117.80	120.30
28	11	236	GLY	N-CA-C	6.25	128.71	113.10
26	14	1786	A	N1-C2-N3	6.25	132.42	129.30
26	14	2374	C	C2-N3-C4	-6.25	116.78	119.90
26	1H	1437	C	C5-C6-N1	6.25	124.12	121.00
26	1H	1978	A	O5'-P-OP2	6.25	118.19	110.70
26	14	1680	U	O5'-P-OP1	-6.25	100.08	105.70
23	2K	40	C	C6-N1-C2	-6.24	117.80	120.30
26	1H	37	C	C4-C5-C6	6.24	120.52	117.40
26	1H	193	U	N1-C2-O2	-6.24	118.43	122.80
26	1H	1142(A)	A	N3-C4-C5	6.24	131.17	126.80
26	1H	2340	G	C8-N9-C4	6.24	108.90	106.40
26	1H	2418	A	C2-N3-C4	6.24	113.72	110.60
26	14	752	A	P-O3'-C3'	6.24	127.19	119.70
26	14	1786	A	N9-C1'-C2'	6.24	122.12	114.00
26	14	2066	C	N3-C4-N4	6.24	122.37	118.00
26	1H	1323	U	N3-C4-C5	-6.24	110.86	114.60
26	1H	1666	G	N3-C4-C5	-6.24	125.48	128.60
26	1H	2070	G	N3-C4-N9	6.24	129.74	126.00
26	1H	996	A	C8-N9-C4	6.24	108.30	105.80
26	14	1982	C	C5-C6-N1	6.24	124.12	121.00
26	1H	58	G	N3-C4-N9	-6.24	122.26	126.00
26	14	1899	G	N7-C8-N9	6.24	116.22	113.10
26	1H	632	A	C5-N7-C8	-6.23	100.78	103.90
26	1H	1636	C	N3-C4-N4	6.23	122.36	118.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
39	A8	24	LEU	CA-CB-CG	6.23	129.64	115.30
26	14	2238	G	O5'-P-OP2	-6.23	100.09	105.70
26	14	2392	A	C5-C6-N1	-6.23	114.58	117.70
1	13	1290	G	N7-C8-N9	6.23	116.22	113.10
26	14	1899	G	C6-C5-N7	-6.23	126.66	130.40
26	14	2445	G	N1-C6-O6	-6.23	116.16	119.90
49	G5	42	GLY	N-CA-C	-6.23	97.52	113.10
23	2K	21	U	C6-N1-C1'	-6.23	112.48	121.20
26	1H	1292	U	O5'-P-OP2	-6.23	100.09	105.70
1	13	748	C	C5-C6-N1	6.23	124.11	121.00
26	1H	1599	C	O5'-P-OP2	-6.23	100.10	105.70
26	14	179	G	C8-N9-C4	6.23	108.89	106.40
26	1H	113	G	C5-C6-N1	-6.22	108.39	111.50
26	14	2591	C	N1-C2-O2	-6.22	115.17	118.90
26	1H	1900	A	OP1-P-OP2	-6.22	110.27	119.60
26	1H	1569	A	C5-N7-C8	-6.22	100.79	103.90
26	1H	239	U	C2-N1-C1'	-6.22	110.24	117.70
26	1H	2461	C	N3-C4-N4	-6.22	113.65	118.00
26	1H	1382	G	OP2-P-O3'	6.22	118.88	105.20
26	14	447	A	N9-C4-C5	6.22	108.29	105.80
26	14	2346	A	N1-C2-N3	6.22	132.41	129.30
1	13	1299	A	N7-C8-N9	6.21	116.91	113.80
26	14	1154	G	C6-C5-N7	-6.21	126.67	130.40
26	14	2067	G	C8-N9-C4	-6.21	103.92	106.40
26	1H	628	G	OP1-P-OP2	6.21	128.92	119.60
26	1H	2018	G	C5-N7-C8	-6.21	101.19	104.30
1	13	1402	C	N3-C4-C5	-6.21	119.42	121.90
26	1H	213	A	C8-N9-C4	6.21	108.28	105.80
26	1H	335	C	N3-C4-C5	-6.21	119.42	121.90
26	1H	796	C	C5-C6-N1	-6.21	117.89	121.00
27	16	6	C	N1-C2-O2	-6.21	115.17	118.90
26	14	457	A	C8-N9-C4	-6.21	103.32	105.80
1	13	1522	U	C4-C5-C6	6.21	123.42	119.70
26	1H	500	G	O5'-P-OP1	-6.21	100.11	105.70
26	1H	692	C	C4-C5-C6	6.21	120.50	117.40
26	1H	974(A)	C	C5-C4-N4	6.21	124.55	120.20
26	14	1776	G	N3-C4-N9	6.21	129.72	126.00
28	19	60	ARG	NE-CZ-NH1	-6.21	117.20	120.30
26	1H	805	G	OP1-P-O3'	6.21	118.86	105.20
26	1H	1698	A	N1-C6-N6	6.21	122.32	118.60
43	E8	51	LEU	CA-CB-CG	6.21	129.58	115.30
26	14	1681	G	N1-C6-O6	6.21	123.62	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	2500	U	C5-C6-N1	-6.21	119.60	122.70
26	14	74	A	C6-C5-N7	-6.21	127.96	132.30
26	1H	690	G	N9-C4-C5	-6.20	102.92	105.40
26	1H	2199	A	O5'-P-OP1	-6.20	100.12	105.70
26	14	1844	C	N3-C2-O2	6.20	126.24	121.90
26	1H	1799	G	P-O3'-C3'	6.20	127.14	119.70
26	1H	1815	A	C6-N1-C2	-6.20	114.88	118.60
26	14	2012	G	C6-C5-N7	-6.20	126.68	130.40
26	14	2821	A	C2-N3-C4	-6.20	107.50	110.60
1	13	266	G	C6-C5-N7	-6.20	126.68	130.40
1	13	767	A	N1-C2-N3	6.20	132.40	129.30
26	1H	1228	G	C2-N3-C4	-6.20	108.80	111.90
26	14	1558	A	C2-N3-C4	-6.20	107.50	110.60
1	13	564	C	C6-N1-C2	-6.20	117.82	120.30
1	13	1053	G	P-O3'-C3'	6.20	127.14	119.70
26	1H	111	A	O5'-P-OP2	-6.20	100.12	105.70
26	1H	1757	U	OP1-P-O3'	6.20	118.83	105.20
39	A8	48	LEU	CA-CB-CG	6.20	129.56	115.30
26	14	1814	G	N9-C4-C5	6.20	107.88	105.40
26	14	2256	G	N1-C2-N2	-6.20	110.62	116.20
26	14	2873	A	C4-N9-C1'	6.20	137.46	126.30
26	1H	1632	A	N1-C6-N6	6.20	122.32	118.60
26	14	979	G	N7-C8-N9	6.20	116.20	113.10
26	14	1400	G	C5-C6-O6	6.20	132.32	128.60
26	14	1619	G	C5-C6-O6	-6.20	124.88	128.60
26	14	2606	C	C2-N3-C4	-6.20	116.80	119.90
38	55	79	LEU	CA-CB-CG	6.20	129.55	115.30
1	13	1227	A	C5-N7-C8	-6.20	100.80	103.90
1	1G	1432	G	C5-C6-N1	-6.20	108.40	111.50
26	14	1950	G	C8-N9-C1'	-6.20	118.95	127.00
26	14	2000	G	C8-N9-C4	6.20	108.88	106.40
26	1H	2374	C	C2-N3-C4	-6.19	116.80	119.90
26	1H	2773	C	C6-N1-C2	6.19	122.78	120.30
1	1G	454	C	N1-C2-O2	6.19	122.62	118.90
26	14	672	C	C2-N1-C1'	-6.19	111.99	118.80
26	1H	651	G	OP1-P-OP2	-6.19	110.31	119.60
26	1H	1617	C	OP2-P-O3'	6.19	118.82	105.20
26	1H	2347	C	OP2-P-O3'	6.19	118.82	105.20
26	14	1283	G	O5'-P-OP2	-6.19	100.13	105.70
26	14	1304	C	N1-C2-O2	6.19	122.61	118.90
26	1H	671	C	C5-C6-N1	-6.19	117.91	121.00
26	1H	930	U	N3-C4-O4	-6.19	115.07	119.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	470	A	N1-C2-N3	6.19	132.40	129.30
26	14	1762	A	C5-C6-N1	-6.19	114.61	117.70
26	14	2553	G	OP1-P-O3'	6.19	118.81	105.20
26	1H	2286	A	N1-C6-N6	6.19	122.31	118.60
26	1H	783	A	N9-C1'-C2'	-6.18	105.20	112.00
26	1H	786	C	OP2-P-O3'	6.18	118.81	105.20
26	1H	1790	C	N1-C2-O2	-6.18	115.19	118.90
26	14	71	A	N1-C2-N3	6.18	132.39	129.30
26	14	791	C	N1-C2-O2	-6.18	115.19	118.90
27	16	61	G	C8-N9-C4	-6.18	103.93	106.40
27	16	98	G	N3-C4-N9	6.18	129.71	126.00
1	1G	913	A	P-O3'-C3'	6.18	127.12	119.70
26	1H	915	C	N1-C2-O2	6.18	122.61	118.90
26	1H	1307	A	C2-N3-C4	-6.18	107.51	110.60
26	1H	2296	U	N3-C4-O4	6.18	123.73	119.40
26	1H	2444	G	N3-C2-N2	-6.18	115.57	119.90
26	14	1319	G	O5'-P-OP2	-6.18	100.14	105.70
26	1H	2449	U	OP2-P-O3'	6.18	118.79	105.20
26	14	2501	C	C6-N1-C1'	6.18	128.21	120.80
1	13	768	A	C6-N1-C2	-6.18	114.89	118.60
26	1H	700	G	N3-C2-N2	-6.18	115.58	119.90
26	1H	1385	G	N3-C4-N9	-6.18	122.29	126.00
26	1H	2582	G	N3-C2-N2	6.18	124.22	119.90
26	14	2442	C	N3-C2-O2	-6.18	117.58	121.90
26	14	2039	C	N3-C4-N4	6.17	122.32	118.00
26	1H	2336	A	O5'-P-OP1	-6.17	100.14	105.70
26	1H	1640	C	OP1-P-OP2	-6.17	110.34	119.60
26	14	330	A	C5-N7-C8	-6.17	100.81	103.90
26	14	386	G	N1-C6-O6	6.17	123.60	119.90
26	1H	1681	G	C5-C6-O6	-6.17	124.90	128.60
26	1H	2713	A	C6-C5-N7	-6.17	127.98	132.30
26	14	137	C	C6-N1-C2	-6.17	117.83	120.30
26	1H	734	A	C6-N1-C2	6.17	122.30	118.60
26	1H	954	G	O5'-P-OP1	-6.17	100.15	105.70
26	1H	2591	C	C5-C4-N4	-6.17	115.88	120.20
26	14	2005	A	N7-C8-N9	-6.17	110.72	113.80
26	1H	391	G	C5-C6-N1	-6.17	108.42	111.50
26	1H	1502	C	C6-N1-C2	-6.16	117.83	120.30
26	1H	271(B)	G	C6-N1-C2	-6.16	121.40	125.10
26	1H	1189	A	N1-C6-N6	6.16	122.30	118.60
26	1H	1363	C	N3-C4-N4	-6.16	113.69	118.00
1	13	789	U	C6-N1-C2	-6.16	117.30	121.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	378	C	C5-C4-N4	-6.16	115.89	120.20
26	1H	378	C	N3-C4-N4	6.16	122.31	118.00
26	1H	1520	U	N1-C2-N3	6.16	118.60	114.90
26	14	2388	A	O4'-C1'-N9	6.16	113.13	108.20
26	1H	906	G	N9-C4-C5	6.16	107.86	105.40
26	14	217	G	O5'-P-OP1	-6.16	100.16	105.70
1	13	129	U	O4'-C1'-N1	6.16	113.12	108.20
26	1H	683	C	C2-N3-C4	-6.16	116.82	119.90
26	1H	1831	G	N7-C8-N9	6.16	116.18	113.10
26	14	2459	A	C8-N9-C4	-6.16	103.34	105.80
26	1H	481	G	O4'-C1'-N9	6.15	113.12	108.20
26	1H	911	A	N1-C6-N6	6.15	122.29	118.60
26	1H	1214	A	OP2-P-O3'	6.15	118.74	105.20
26	14	2019	A	N7-C8-N9	-6.15	110.72	113.80
26	14	2039	C	C6-N1-C2	-6.15	117.84	120.30
26	14	133	C	C2-N3-C4	-6.15	116.83	119.90
26	1H	263	C	N1-C2-O2	6.15	122.59	118.90
26	1H	793	A	C6-N1-C2	-6.15	114.91	118.60
26	1H	1161	C	C6-N1-C2	-6.15	117.84	120.30
26	14	2584	U	C6-N1-C1'	-6.15	112.59	121.20
26	14	1661	G	N9-C4-C5	-6.15	102.94	105.40
26	1H	1936	A	C4-C5-N7	6.14	113.77	110.70
27	16	98	G	N3-C2-N2	6.14	124.20	119.90
26	14	71	A	N7-C8-N9	6.14	116.87	113.80
26	1H	1161	C	C5-C6-N1	6.14	124.07	121.00
26	1H	2380	C	C4-C5-C6	6.14	120.47	117.40
1	1G	719	C	C6-N1-C2	-6.14	117.84	120.30
26	14	2258	C	N1-C2-O2	-6.14	115.22	118.90
26	1H	1623	G	C5-C6-O6	6.14	132.28	128.60
26	14	119	A	N1-C2-N3	6.14	132.37	129.30
26	1H	853	G	O5'-P-OP1	6.14	118.07	110.70
26	14	143	C	C6-N1-C2	-6.14	117.84	120.30
26	1H	624	C	N3-C2-O2	6.14	126.20	121.90
26	1H	2378	A	N1-C6-N6	6.14	122.28	118.60
26	14	1021	A	C2-N3-C4	-6.14	107.53	110.60
26	14	2873	A	C4-C5-C6	6.14	120.07	117.00
5	4E	91	LEU	CA-CB-CG	6.13	129.41	115.30
26	1H	1782	C	C2-N3-C4	-6.13	116.83	119.90
1	1G	115	G	N1-C6-O6	-6.13	116.22	119.90
26	14	915	C	N1-C2-O2	6.13	122.58	118.90
26	14	1333	C	C5-C4-N4	-6.13	115.91	120.20
26	14	1698	A	N1-C2-N3	6.13	132.37	129.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	2258	C	N3-C2-O2	6.13	126.19	121.90
26	1H	227	A	N1-C6-N6	-6.13	114.92	118.60
26	1H	609	A	N1-C6-N6	6.13	122.28	118.60
1	13	1203	C	C6-N1-C2	-6.13	117.85	120.30
26	1H	784	A	N9-C4-C5	6.13	108.25	105.80
26	1H	1191	G	C8-N9-C4	6.13	108.85	106.40
26	1H	2394	C	O5'-P-OP2	-6.13	100.18	105.70
26	14	1786	A	N1-C6-N6	6.13	122.28	118.60
26	1H	265	A	C6-C5-N7	-6.13	128.01	132.30
26	1H	624	C	N3-C4-N4	6.13	122.29	118.00
26	1H	988	A	N7-C8-N9	6.13	116.86	113.80
26	1H	1754	C	C6-N1-C2	-6.13	117.85	120.30
26	1H	2430	A	C5-C6-N6	-6.13	118.80	123.70
26	1H	1228	G	N1-C2-N3	6.13	127.58	123.90
1	13	1329	A	N1-C6-N6	6.13	122.28	118.60
26	1H	596	G	N3-C2-N2	-6.13	115.61	119.90
26	1H	2300	G	C8-N9-C4	-6.13	103.95	106.40
26	1H	2837	G	N1-C6-O6	6.13	123.58	119.90
26	1H	2856	C	C6-N1-C2	-6.13	117.85	120.30
26	14	2329	G	N1-C6-O6	-6.13	116.22	119.90
26	1H	764	A	OP1-P-OP2	-6.12	110.41	119.60
26	1H	1613	G	N3-C4-N9	6.12	129.68	126.00
26	1H	1899	G	C5-N7-C8	-6.12	101.24	104.30
1	1G	904	C	O5'-P-OP1	-6.12	100.19	105.70
26	14	2439	A	C6-C5-N7	-6.12	128.01	132.30
26	1H	62	C	C5-C6-N1	-6.12	117.94	121.00
26	1H	831	G	C8-N9-C4	6.12	108.85	106.40
26	1H	1379	A	C6-C5-N7	-6.12	128.01	132.30
26	14	189	G	C8-N9-C4	6.12	108.85	106.40
26	14	803	U	O5'-P-OP1	6.12	118.04	110.70
26	14	2700	C	C6-N1-C1'	-6.12	113.46	120.80
26	1H	2318	G	N7-C8-N9	6.11	116.16	113.10
26	14	195	A	P-O3'-C3'	6.11	127.04	119.70
26	14	2031	A	N1-C6-N6	6.11	122.27	118.60
22	1K	76	A	O4'-C1'-N9	6.11	113.09	108.20
26	14	672	C	O5'-P-OP2	-6.11	100.20	105.70
26	14	1334	G	N7-C8-N9	6.11	116.16	113.10
26	14	1489	U	C5-C4-O4	6.11	129.57	125.90
1	13	112	G	C8-N9-C4	-6.11	103.96	106.40
1	13	137	C	C6-N1-C2	6.11	122.74	120.30
26	1H	1010	A	C8-N9-C4	6.11	108.24	105.80
26	1H	2752	C	C5-C6-N1	6.11	124.05	121.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	71	A	O4'-C1'-N9	-6.11	103.31	108.20
26	1H	207	A	C2-N3-C4	-6.11	107.55	110.60
26	1H	1998	G	N7-C8-N9	-6.11	110.05	113.10
26	1H	2392	A	N3-C4-C5	6.11	131.07	126.80
23	2L	48	U	P-O3'-C3'	6.10	127.03	119.70
1	13	903	G	O5'-P-OP2	-6.10	100.21	105.70
26	14	1785	A	C4-C5-C6	6.10	120.05	117.00
26	14	2031	A	C5-C6-N6	-6.10	118.82	123.70
26	1H	1187	G	OP2-P-O3'	6.10	118.62	105.20
1	13	728	A	C8-N9-C4	-6.10	103.36	105.80
26	1H	641	C	O5'-P-OP2	6.10	118.02	110.70
26	14	747	U	N3-C2-O2	6.10	126.47	122.20
26	14	2543	G	C8-N9-C4	6.10	108.84	106.40
26	1H	1596	A	N1-C6-N6	-6.10	114.94	118.60
26	1H	2299	G	C5-N7-C8	-6.10	101.25	104.30
26	1H	2375	G	N9-C4-C5	-6.10	102.96	105.40
26	14	1826	G	C5-N7-C8	6.10	107.35	104.30
26	1H	1789	A	C5-C6-N1	6.10	120.75	117.70
26	1H	1981	A	OP2-P-O3'	6.10	118.61	105.20
26	14	2712	U	N3-C4-O4	-6.10	115.13	119.40
26	1H	839	U	C4-C5-C6	6.09	123.36	119.70
26	1H	913	U	OP1-P-OP2	6.09	128.74	119.60
26	1H	2249	U	N3-C4-C5	-6.09	110.94	114.60
26	14	1613	G	C5-N7-C8	6.09	107.35	104.30
26	14	2501	C	N3-C4-C5	6.09	124.34	121.90
26	1H	1231	G	C5-C6-O6	-6.09	124.94	128.60
26	1H	1314	C	C6-N1-C1'	-6.09	113.49	120.80
26	1H	2329	G	OP1-P-OP2	6.09	128.74	119.60
26	14	864	G	OP1-P-OP2	-6.09	110.47	119.60
26	1H	127	A	C4-C5-N7	6.09	113.74	110.70
26	14	1517	G	OP1-P-O3'	6.09	118.59	105.20
26	1H	329	G	O5'-P-OP2	-6.09	100.22	105.70
26	14	831	G	C8-N9-C4	6.09	108.83	106.40
26	14	2373	G	N3-C2-N2	-6.09	115.64	119.90
26	1H	1899	G	C8-N9-C1'	6.08	134.91	127.00
26	1H	2281	C	N3-C4-N4	6.08	122.26	118.00
26	14	1636	C	N3-C4-N4	6.08	122.26	118.00
26	1H	579	G	N1-C2-N2	6.08	121.67	116.20
26	1H	1255	U	N3-C4-C5	-6.08	110.95	114.60
26	1H	1587	A	C8-N9-C4	-6.08	103.37	105.80
26	14	330	A	C5-C6-N1	-6.08	114.66	117.70
26	1H	452	G	C2-N3-C4	6.08	114.94	111.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	1G	1356	G	C8-N9-C4	-6.08	103.97	106.40
26	14	1767	C	C5-C6-N1	-6.08	117.96	121.00
26	14	2078	C	C6-N1-C2	-6.08	117.87	120.30
1	13	516	U	C6-N1-C2	-6.08	117.35	121.00
1	13	967	C	N3-C2-O2	6.08	126.15	121.90
26	1H	1187	G	N1-C6-O6	6.08	123.55	119.90
26	1H	2331	G	C2-N3-C4	-6.08	108.86	111.90
1	13	413	G	O4'-C1'-N9	6.07	113.06	108.20
26	14	470	A	O5'-P-OP1	-6.07	100.23	105.70
26	14	1790	C	OP1-P-O3'	6.07	118.56	105.20
1	13	810	C	O5'-P-OP2	-6.07	100.23	105.70
26	1H	2337	G	C5-C6-O6	-6.07	124.96	128.60
26	14	2250	G	C8-N9-C4	-6.07	103.97	106.40
26	1H	501	A	O5'-P-OP2	-6.07	100.24	105.70
1	1G	1126	U	P-O3'-C3'	6.07	126.98	119.70
26	1H	570	G	C5-C6-N1	-6.07	108.47	111.50
26	1H	951	C	N3-C4-C5	6.07	124.33	121.90
26	14	2001	A	C8-N9-C4	6.07	108.23	105.80
1	13	1158	C	C6-N1-C2	-6.07	117.87	120.30
26	1H	464	U	C4-C5-C6	6.07	123.34	119.70
26	1H	762	U	C5-C6-N1	6.07	125.73	122.70
26	1H	957	A	C4-C5-C6	6.07	120.03	117.00
26	14	1901	A	OP1-P-O3'	6.07	118.55	105.20
26	1H	1674	G	O4'-C1'-N9	-6.07	103.35	108.20
26	1H	2699	C	C5-C4-N4	-6.07	115.95	120.20
26	1H	976	C	O5'-P-OP1	-6.06	100.24	105.70
26	1H	2293	C	N3-C2-O2	-6.06	117.66	121.90
26	1H	1616	A	C6-C5-N7	-6.06	128.06	132.30
26	1H	1915	U	C2-N3-C4	-6.06	123.36	127.00
26	1H	2006	C	O5'-P-OP1	6.06	117.97	110.70
26	1H	2282	G	C5-C6-O6	6.06	132.24	128.60
26	14	2321	G	N3-C4-C5	-6.06	125.57	128.60
26	1H	2422	A	N1-C6-N6	-6.06	114.97	118.60
26	1H	1204	A	C8-N9-C1'	-6.06	116.80	127.70
26	1H	1318	C	O5'-P-OP1	-6.06	100.25	105.70
26	1H	1612	C	C4-C5-C6	6.06	120.43	117.40
26	1H	2638	G	N3-C2-N2	6.06	124.14	119.90
1	1G	267	C	O5'-P-OP1	-6.06	100.25	105.70
1	1G	613	C	C5-C6-N1	6.06	124.03	121.00
26	14	835	A	C2-N3-C4	6.06	113.63	110.60
26	1H	735	A	N9-C4-C5	-6.06	103.38	105.80
26	14	1598	C	OP1-P-OP2	-6.06	110.52	119.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	2389	G	C8-N9-C4	-6.06	103.98	106.40
26	14	1379	A	C5-N7-C8	-6.05	100.87	103.90
26	14	1839	G	O5'-P-OP1	-6.05	100.25	105.70
26	1H	2080	G	O4'-C1'-N9	-6.05	103.36	108.20
26	1H	2244	U	O5'-P-OP2	6.05	117.96	110.70
1	1G	866	C	C6-N1-C2	-6.05	117.88	120.30
26	1H	1275	A	N1-C6-N6	6.05	122.23	118.60
26	1H	1396	U	O5'-P-OP1	-6.05	100.25	105.70
26	1H	2406	U	N1-C2-O2	6.05	127.04	122.80
26	1H	2508	G	N3-C4-N9	-6.05	122.37	126.00
26	14	1618	A	N7-C8-N9	6.05	116.83	113.80
26	1H	1162	G	C8-N9-C4	-6.05	103.98	106.40
24	3L	71	G	N1-C6-O6	-6.05	116.27	119.90
26	14	558	G	N7-C8-N9	-6.05	110.08	113.10
26	14	1965	C	C6-N1-C2	6.05	122.72	120.30
26	1H	1166	C	C5-C6-N1	6.05	124.02	121.00
1	13	805	C	C6-N1-C2	-6.05	117.88	120.30
26	1H	1265	A	C5'-C4'-C3'	-6.05	106.32	116.00
26	1H	2731	G	N1-C6-O6	6.05	123.53	119.90
1	1G	413	G	C6-C5-N7	6.05	134.03	130.40
1	1G	1498	U	P-O3'-C3'	6.05	126.96	119.70
26	14	695	G	N1-C6-O6	6.05	123.53	119.90
26	1H	635	C	O5'-P-OP1	6.04	117.95	110.70
26	1H	1673	U	C2-N1-C1'	-6.04	110.45	117.70
26	1H	2248	C	C6-N1-C2	-6.04	117.88	120.30
1	1G	317	G	C6-C5-N7	-6.04	126.77	130.40
26	14	1252	G	O4'-C1'-N9	-6.04	103.37	108.20
1	13	827	U	C5-C4-O4	6.04	129.53	125.90
26	1H	778	G	C5-N7-C8	6.04	107.32	104.30
26	14	469	G	C5-C6-N1	6.04	114.52	111.50
1	13	1299	A	C5-N7-C8	-6.04	100.88	103.90
26	1H	238	C	C5-C6-N1	-6.04	117.98	121.00
26	1H	763	G	N1-C6-O6	6.04	123.52	119.90
26	1H	1778	U	OP2-P-O3'	6.04	118.49	105.20
26	1H	1905	C	P-O3'-C3'	6.04	126.95	119.70
26	1H	2689	U	N3-C4-O4	-6.04	115.17	119.40
26	14	133	C	C5-C6-N1	-6.04	117.98	121.00
26	1H	478	A	N1-C2-N3	6.04	132.32	129.30
26	1H	842	G	C4-N9-C1'	-6.04	118.65	126.50
1	1G	690	G	C2-N3-C4	-6.04	108.88	111.90
26	14	1787	A	O5'-P-OP1	-6.04	100.27	105.70
26	14	1953	A	O5'-P-OP2	6.04	117.94	110.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	1307	A	OP1-P-OP2	6.04	128.65	119.60
26	1H	1313	U	C6-N1-C2	-6.04	117.38	121.00
26	1H	1446	C	C6-N1-C2	-6.04	117.89	120.30
26	1H	2295	C	N3-C4-C5	-6.04	119.49	121.90
26	14	1332	G	N1-C2-N3	6.04	127.52	123.90
26	1H	1368	G	O5'-P-OP2	-6.03	100.27	105.70
26	14	1804	C	C5-C6-N1	6.03	124.02	121.00
26	14	2546	U	OP1-P-OP2	6.03	128.65	119.60
26	1H	529	A	C5-N7-C8	-6.03	100.89	103.90
26	1H	2275	C	O5'-P-OP2	-6.03	100.27	105.70
26	14	2581	G	N7-C8-N9	6.03	116.12	113.10
1	13	266	G	C4-C5-N7	6.03	113.21	110.80
1	13	956	U	N3-C4-C5	-6.03	110.98	114.60
26	1H	2070	G	N7-C8-N9	-6.03	110.08	113.10
26	1H	1298	C	OP2-P-O3'	-6.03	91.94	105.20
27	16	92	G	OP2-P-O3'	6.03	118.46	105.20
26	14	621	A	N3-C4-C5	6.03	131.02	126.80
26	14	921	G	N7-C8-N9	6.03	116.11	113.10
1	13	449	C	N3-C2-O2	-6.03	117.68	121.90
27	16	13	A	OP1-P-OP2	6.03	128.64	119.60
1	13	268	C	O5'-P-OP2	6.02	117.93	110.70
53	O8	42	TRP	CA-CB-CG	6.02	125.14	113.70
26	14	508	G	O5'-P-OP1	-6.02	100.28	105.70
26	1H	1356	G	O5'-P-OP1	-6.02	100.28	105.70
26	1H	1513	C	C5-C6-N1	6.02	124.01	121.00
26	1H	2436	G	N3-C2-N2	-6.02	115.68	119.90
26	1H	2287	A	N1-C6-N6	6.02	122.21	118.60
26	14	1302	A	N1-C6-N6	-6.02	114.99	118.60
1	13	757	U	N1-C2-O2	6.02	127.01	122.80
26	1H	869	G	N1-C2-N2	-6.02	110.78	116.20
26	1H	1197	G	N7-C8-N9	-6.02	110.09	113.10
26	1H	1694	C	P-O3'-C3'	6.02	126.92	119.70
26	1H	2688	U	C6-N1-C2	-6.02	117.39	121.00
1	13	812	C	N3-C2-O2	-6.02	117.69	121.90
1	13	827	U	C4-C5-C6	6.02	123.31	119.70
26	1H	501	A	OP1-P-OP2	6.02	128.63	119.60
26	1H	598	G	OP1-P-OP2	6.02	128.63	119.60
22	1K	18	G	C4-N9-C1'	6.02	134.32	126.50
26	1H	1969	A	C5-C6-N6	6.02	128.51	123.70
26	1H	2070	G	C5-N7-C8	6.02	107.31	104.30
26	1H	2366	A	OP2-P-O3'	6.01	118.43	105.20
26	1H	2440	C	N3-C4-C5	-6.01	119.49	121.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	1334	G	O5'-P-OP2	6.01	117.92	110.70
26	14	2281	C	C6-N1-C2	-6.01	117.89	120.30
26	14	140	A	C6-C5-N7	-6.01	128.09	132.30
26	14	386	G	C4-C5-N7	6.01	113.20	110.80
26	1H	74	A	O4'-C1'-N9	-6.01	103.39	108.20
26	1H	863	A	N7-C8-N9	-6.01	110.79	113.80
1	1G	353	A	C4-C5-N7	6.01	113.71	110.70
26	14	89	G	N1-C6-O6	6.01	123.51	119.90
26	1H	1308	A	N9-C4-C5	6.01	108.20	105.80
26	1H	1616	A	C2-N3-C4	-6.01	107.60	110.60
26	1H	2690	C	C2-N1-C1'	-6.01	112.19	118.80
26	14	1258	C	C4-C5-C6	-6.01	114.39	117.40
1	13	1222	G	O5'-P-OP2	-6.01	100.29	105.70
26	1H	201	C	C6-N1-C2	6.01	122.70	120.30
26	1H	1639	U	N1-C2-O2	6.01	127.00	122.80
26	14	827	U	O5'-P-OP2	-6.01	100.29	105.70
26	1H	71	A	C8-N9-C4	-6.01	103.40	105.80
26	1H	1271	G	N1-C6-O6	6.01	123.50	119.90
26	1H	1900	A	N9-C4-C5	-6.01	103.40	105.80
26	1H	2718	G	N1-C6-O6	6.01	123.50	119.90
26	14	245	G	C5-C6-O6	-6.01	125.00	128.60
26	14	574	C	C2-N1-C1'	-6.01	112.19	118.80
2	12	196	LEU	CA-CB-CG	6.00	129.11	115.30
26	1H	1201	C	N1-C2-O2	-6.00	115.30	118.90
26	1H	1347	G	OP1-P-O3'	6.00	118.41	105.20
26	1H	1489	U	C5-C4-O4	6.00	129.50	125.90
26	1H	1982	C	C5-C6-N1	6.00	124.00	121.00
26	1H	2822	G	C5-C6-O6	-6.00	125.00	128.60
26	1H	463	G	O5'-P-OP1	6.00	117.90	110.70
26	14	565	C	N1-C2-O2	6.00	122.50	118.90
26	14	2061	G	C8-N9-C4	6.00	108.80	106.40
1	1G	266	G	P-O3'-C3'	6.00	126.90	119.70
30	39	125	LEU	CA-CB-CG	6.00	129.09	115.30
23	2K	35	C	OP1-P-O3'	6.00	118.39	105.20
26	1H	1307	A	O5'-P-OP1	-6.00	100.30	105.70
1	1G	218	C	C6-N1-C2	-6.00	117.90	120.30
24	3L	71	G	O4'-C1'-N9	6.00	113.00	108.20
26	14	465	G	N3-C4-N9	6.00	129.60	126.00
26	1H	1259	G	OP2-P-O3'	6.00	118.39	105.20
26	1H	566	U	C6-N1-C2	5.99	124.60	121.00
26	14	2558	C	C5-C4-N4	-5.99	116.00	120.20
26	1H	391	G	C4-C5-C6	5.99	122.39	118.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	871	U	O5'-P-OP1	-5.99	100.31	105.70
26	14	1380	G	C2-N3-C4	-5.99	108.90	111.90
26	1H	1900	A	N3-C4-N9	5.99	132.19	127.40
26	1H	382	G	OP1-P-O3'	5.99	118.38	105.20
26	1H	803	U	C5-C6-N1	-5.99	119.71	122.70
1	1G	1514	C	C6-N1-C2	-5.99	117.90	120.30
26	14	565	C	C5-C6-N1	-5.99	118.01	121.00
26	1H	189	G	C5-C6-O6	-5.99	125.01	128.60
26	14	201	C	C2-N3-C4	-5.99	116.91	119.90
26	14	1239	G	N1-C6-O6	5.99	123.49	119.90
26	14	2606	C	O5'-P-OP1	-5.99	100.31	105.70
26	1H	940	G	N3-C4-C5	-5.98	125.61	128.60
1	1G	1157	A	P-O3'-C3'	5.98	126.88	119.70
1	13	1459	C	N1-C2-O2	5.98	122.49	118.90
26	1H	1967	C	OP1-P-OP2	5.98	128.57	119.60
26	1H	2705	A	C6-C5-N7	-5.98	128.11	132.30
26	14	737	C	C5-C4-N4	-5.98	116.01	120.20
26	14	1348	G	N1-C6-O6	5.98	123.49	119.90
1	13	690	G	C5-C6-N1	-5.98	108.51	111.50
26	1H	35	G	O5'-P-OP2	-5.98	100.32	105.70
26	1H	2507	C	N3-C4-C5	-5.98	119.51	121.90
1	1G	581	G	O5'-P-OP2	-5.98	100.32	105.70
25	4K	19	U	O5'-P-OP2	-5.98	100.32	105.70
1	1G	932	C	C2-N1-C1'	5.98	125.38	118.80
26	14	1332	G	C8-N9-C4	-5.98	104.01	106.40
26	14	1816	G	C5-C6-O6	-5.98	125.01	128.60
1	13	956	U	C5-C6-N1	5.98	125.69	122.70
26	1H	682	G	O5'-P-OP2	-5.98	100.32	105.70
26	1H	1398	C	OP1-P-OP2	-5.98	110.63	119.60
1	13	690	G	C4-C5-N7	5.98	113.19	110.80
1	13	1432	G	C5-C6-O6	-5.97	125.02	128.60
26	1H	134	C	C2-N3-C4	-5.97	116.91	119.90
26	1H	1611	C	C6-N1-C2	5.97	122.69	120.30
26	1H	1969	A	C4-C5-N7	-5.97	107.71	110.70
1	13	975	A	O4'-C1'-N9	-5.97	103.42	108.20
26	1H	2385	C	C2-N3-C4	-5.97	116.91	119.90
26	14	208	C	N3-C4-N4	5.97	122.18	118.00
26	14	1248	G	OP1-P-OP2	-5.97	110.64	119.60
26	14	1248	G	O5'-P-OP1	5.97	117.87	110.70
1	13	1201	A	N1-C6-N6	5.97	122.18	118.60
26	1H	2387	U	OP2-P-O3'	5.97	118.34	105.20
1	13	1025	U	C5-C6-N1	5.97	125.69	122.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	835	A	C2-N3-C4	5.97	113.58	110.60
26	14	570	G	C8-N9-C1'	-5.97	119.24	127.00
26	14	755	C	N1-C2-O2	-5.97	115.32	118.90
26	14	1698	A	N9-C4-C5	-5.97	103.41	105.80
26	1H	139	G	C2-N3-C4	5.96	114.88	111.90
26	1H	1559	G	N3-C4-N9	-5.96	122.42	126.00
26	14	830	G	N9-C4-C5	-5.96	103.01	105.40
26	14	911	A	OP1-P-O3'	5.96	118.32	105.20
26	1H	198	C	N3-C4-C5	5.96	124.28	121.90
26	14	2581	G	N3-C4-C5	-5.96	125.62	128.60
1	13	1279	A	C8-N9-C4	-5.96	103.42	105.80
26	1H	2286	A	C6-C5-N7	-5.96	128.13	132.30
1	1G	687	A	P-O3'-C3'	5.96	126.85	119.70
26	14	677	A	OP1-P-OP2	5.96	128.54	119.60
26	14	1367	A	C6-C5-N7	-5.96	128.13	132.30
26	14	2503	A	O5'-P-OP2	-5.96	100.33	105.70
26	1H	1141	U	N3-C2-O2	-5.96	118.03	122.20
26	1H	1515	C	C6-N1-C2	-5.96	117.92	120.30
26	14	2365	G	C5-C6-O6	-5.96	125.03	128.60
26	1H	121	G	N9-C4-C5	-5.96	103.02	105.40
26	1H	1941	C	N3-C4-C5	-5.96	119.52	121.90
26	1H	2312	U	O5'-P-OP1	-5.96	100.34	105.70
26	1H	2562	U	C5-C6-N1	-5.96	119.72	122.70
26	14	2049	G	N3-C2-N2	-5.96	115.73	119.90
1	13	1432	G	N1-C6-O6	5.95	123.47	119.90
26	1H	407	G	N3-C4-C5	-5.95	125.62	128.60
27	16	29	A	C8-N9-C4	-5.95	103.42	105.80
26	14	2463	C	N1-C2-O2	-5.95	115.33	118.90
26	14	2707	G	O4'-C1'-N9	5.95	112.96	108.20
1	13	977	A	C8-N9-C4	-5.95	103.42	105.80
23	2K	13	C	C2-N3-C4	5.95	122.88	119.90
26	1H	805	G	C4-C5-C6	5.95	122.37	118.80
26	1H	920	G	C8-N9-C4	5.95	108.78	106.40
26	1H	1653	G	N3-C4-N9	5.95	129.57	126.00
26	1H	2031	A	N1-C2-N3	-5.95	126.32	129.30
26	14	1418	G	C6-C5-N7	-5.95	126.83	130.40
26	14	2426	A	OP1-P-O3'	5.95	118.29	105.20
26	1H	1241	A	C5-N7-C8	-5.95	100.93	103.90
26	1H	2299	G	N1-C6-O6	5.95	123.47	119.90
1	13	525	C	C5-C4-N4	-5.95	116.04	120.20
26	1H	13	A	C8-N9-C4	-5.95	103.42	105.80
26	1H	906	G	C6-C5-N7	5.95	133.97	130.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	1193	G	O5'-P-OP1	5.95	117.84	110.70
1	13	115	G	P-O3'-C3'	5.95	126.83	119.70
26	1H	1224	G	C4-N9-C1'	-5.95	118.77	126.50
1	13	967	C	N1-C2-O2	-5.94	115.33	118.90
26	1H	1349	A	N9-C4-C5	-5.94	103.42	105.80
26	14	1930	G	C4-C5-N7	-5.94	108.42	110.80
26	1H	1298	C	N1-C2-O2	5.94	122.47	118.90
26	1H	2248	C	C5-C4-N4	5.94	124.36	120.20
1	13	1336	C	P-O3'-C3'	5.94	126.83	119.70
26	1H	2273	A	N1-C2-N3	-5.94	126.33	129.30
26	1H	2546	U	N3-C2-O2	5.94	126.36	122.20
26	14	1972	A	OP2-P-O3'	5.94	118.27	105.20
26	14	2004	G	N1-C6-O6	5.94	123.46	119.90
26	1H	459	U	N1-C2-N3	5.94	118.46	114.90
26	1H	1785	A	C8-N9-C4	-5.94	103.42	105.80
26	1H	2465	C	C5-C6-N1	-5.94	118.03	121.00
26	1H	2336	A	C2-N3-C4	5.94	113.57	110.60
26	1H	2819	G	C6-C5-N7	-5.94	126.84	130.40
26	1H	793	A	N3-C4-C5	-5.93	122.64	126.80
26	1H	2576	G	N9-C4-C5	-5.93	103.03	105.40
26	1H	2591	C	N3-C4-N4	5.93	122.15	118.00
26	1H	2686	G	N3-C4-N9	5.93	129.56	126.00
26	14	856	C	C2-N1-C1'	5.93	125.33	118.80
26	14	1145	C	C5-C6-N1	5.93	123.97	121.00
1	1G	723	U	C5-C6-N1	5.93	125.67	122.70
26	1H	2006	C	C6-N1-C2	5.93	122.67	120.30
26	14	771	G	OP1-P-O3'	5.93	118.25	105.20
26	1H	2584	U	N1-C2-N3	5.93	118.46	114.90
26	14	123	G	C5-C6-O6	-5.93	125.04	128.60
26	1H	71	A	N1-C6-N6	5.93	122.16	118.60
26	1H	536	A	N9-C4-C5	5.93	108.17	105.80
26	1H	1418	G	N1-C6-O6	-5.93	116.34	119.90
26	1H	2016	U	N3-C4-O4	-5.93	115.25	119.40
27	16	81	G	N7-C8-N9	5.93	116.06	113.10
1	1G	353	A	C5-N7-C8	-5.93	100.94	103.90
26	14	794	G	C4-C5-N7	-5.93	108.43	110.80
26	1H	51	G	O4'-C1'-N9	-5.92	103.46	108.20
26	1H	1888	G	C4-N9-C1'	5.92	134.20	126.50
26	14	787	U	O5'-P-OP1	5.92	117.81	110.70
26	14	1302	A	OP1-P-OP2	5.92	128.49	119.60
26	14	1615	C	N3-C4-C5	-5.92	119.53	121.90
1	13	1519	A	C8-N9-C4	-5.92	103.43	105.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	372	G	N9-C4-C5	5.92	107.77	105.40
26	1H	755	C	N3-C4-C5	-5.92	119.53	121.90
26	1H	1348	G	O5'-P-OP2	5.92	117.81	110.70
22	1K	18	G	P-O3'-C3'	5.92	126.81	119.70
26	1H	775	G	O4'-C1'-N9	5.92	112.94	108.20
26	1H	1669	A	N7-C8-N9	5.92	116.76	113.80
26	1H	2243	U	OP2-P-O3'	5.92	118.22	105.20
26	1H	2375	G	C5-C6-N1	5.92	114.46	111.50
26	14	1309	G	N9-C4-C5	-5.92	103.03	105.40
26	1H	622	G	N7-C8-N9	-5.92	110.14	113.10
26	1H	914	C	C2-N1-C1'	-5.92	112.29	118.80
1	1G	664	G	N3-C4-N9	-5.92	122.45	126.00
1	1G	772	U	O5'-P-OP2	-5.92	100.37	105.70
26	14	138	G	C8-N9-C4	-5.92	104.03	106.40
26	1H	180	G	N9-C4-C5	-5.92	103.03	105.40
26	1H	808	G	N1-C2-N2	-5.92	110.88	116.20
26	1H	1790	C	N3-C4-C5	5.92	124.27	121.90
26	14	1141	U	P-O3'-C3'	5.92	126.80	119.70
26	14	1380	G	C5-C6-N1	-5.92	108.54	111.50
26	1H	663	G	N9-C4-C5	5.91	107.77	105.40
26	1H	1614	A	N3-C4-N9	-5.91	122.67	127.40
26	1H	2074	U	C5-C6-N1	-5.91	119.74	122.70
26	1H	2289	G	N1-C2-N2	5.91	121.52	116.20
26	14	857	C	C6-N1-C2	-5.91	117.93	120.30
26	14	1274	A	C4-C5-C6	5.91	119.96	117.00
26	14	2571	C	C5-C4-N4	-5.91	116.06	120.20
1	1G	1529	G	C8-N9-C4	-5.91	104.03	106.40
26	1H	113	G	C2-N3-C4	-5.91	108.94	111.90
26	1H	974(A)	C	N3-C4-C5	-5.91	119.54	121.90
26	1H	1340	U	C5-C4-O4	-5.91	122.35	125.90
26	1H	2065	C	N3-C2-O2	-5.91	117.76	121.90
1	13	797	C	O5'-P-OP1	-5.91	100.38	105.70
26	1H	354	G	N1-C6-O6	5.91	123.45	119.90
26	14	945	A	N7-C8-N9	5.91	116.75	113.80
26	14	1376	C	O5'-P-OP1	-5.91	100.38	105.70
26	14	1574	C	OP2-P-O3'	5.91	118.20	105.20
1	13	1299	A	C6-C5-N7	-5.91	128.16	132.30
27	16	60	C	C6-N1-C2	-5.91	117.94	120.30
26	14	2265	U	N3-C4-O4	5.91	123.53	119.40
26	1H	321	G	N1-C6-O6	5.91	123.44	119.90
26	1H	1197	G	C8-N9-C4	5.91	108.76	106.40
26	1H	2688	U	N3-C4-O4	-5.91	115.27	119.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	1G	1408	A	N1-C6-N6	-5.91	115.06	118.60
26	14	265	A	C2-N3-C4	-5.91	107.65	110.60
26	14	1798	U	O5'-P-OP2	-5.91	100.39	105.70
26	14	2392	A	C2-N3-C4	-5.91	107.65	110.60
26	1H	517	C	N3-C4-N4	5.90	122.13	118.00
26	1H	209	C	O5'-P-OP2	-5.90	100.39	105.70
26	1H	2398	U	N1-C2-O2	5.90	126.93	122.80
24	3L	71	G	C5-C6-O6	5.90	132.14	128.60
26	14	121	G	C6-N1-C2	-5.90	121.56	125.10
26	14	1971	A	OP2-P-O3'	5.90	118.19	105.20
26	1H	606	U	O5'-P-OP2	-5.90	100.39	105.70
26	1H	735	A	N1-C6-N6	5.90	122.14	118.60
26	1H	802	A	OP1-P-O3'	-5.90	92.22	105.20
26	1H	908	C	OP2-P-O3'	5.90	118.18	105.20
26	1H	2249	U	C5-C6-N1	5.90	125.65	122.70
26	1H	2270	G	N1-C6-O6	5.90	123.44	119.90
26	1H	2680	C	N1-C2-O2	-5.90	115.36	118.90
26	14	215	G	C5-C6-O6	-5.90	125.06	128.60
26	14	2356	C	C2-N1-C1'	-5.90	112.31	118.80
26	1H	928	G	N1-C6-O6	5.90	123.44	119.90
26	1H	2747	G	N3-C4-N9	5.90	129.54	126.00
26	1H	1761	C	C5-C4-N4	-5.90	116.07	120.20
26	1H	2686	G	N3-C4-C5	-5.90	125.65	128.60
26	14	1769	G	N1-C6-O6	5.90	123.44	119.90
26	14	2623	G	O5'-P-OP2	-5.90	100.39	105.70
26	14	1162	G	O5'-P-OP1	-5.90	100.39	105.70
26	14	1614	A	N1-C6-N6	5.90	122.14	118.60
1	13	1103	C	C6-N1-C2	-5.89	117.94	120.30
26	1H	1702	G	C8-N9-C4	5.89	108.76	106.40
26	14	946	G	N9-C4-C5	-5.89	103.04	105.40
26	14	998	C	N1-C2-O2	5.89	122.44	118.90
26	14	2075	U	C2-N3-C4	-5.89	123.46	127.00
1	13	901	A	C2-N3-C4	-5.89	107.65	110.60
26	1H	1761	C	N3-C2-O2	5.89	126.03	121.90
26	14	2681	C	C5-C4-N4	5.89	124.33	120.20
26	1H	513	A	C8-N9-C4	-5.89	103.44	105.80
26	1H	1223	C	N3-C2-O2	5.89	126.02	121.90
26	1H	1784	A	N1-C2-N3	5.89	132.24	129.30
24	3L	71	G	C4-C5-N7	-5.89	108.44	110.80
1	13	428	G	C8-N9-C4	-5.89	104.05	106.40
1	1G	1474	G	N3-C4-C5	5.89	131.54	128.60
26	14	2489	G	OP2-P-O3'	5.89	118.15	105.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	791	C	OP2-P-O3'	5.88	118.15	105.20
26	1H	1900	A	C2-N3-C4	5.88	113.54	110.60
26	1H	2335	A	O4'-C1'-N9	5.88	112.91	108.20
26	1H	837	C	C5-C6-N1	5.88	123.94	121.00
26	1H	866	A	C4-C5-N7	5.88	113.64	110.70
26	1H	1817	G	N9-C4-C5	-5.88	103.05	105.40
26	1H	2396	G	C8-N9-C1'	5.88	134.65	127.00
1	1G	326	G	C5-C6-N1	-5.88	108.56	111.50
26	14	1804	C	C6-N1-C2	-5.88	117.95	120.30
26	1H	1649	G	N3-C4-C5	-5.88	125.66	128.60
26	1H	2392	A	C8-N9-C4	-5.88	103.45	105.80
26	1H	2509	G	C5-C6-N1	5.88	114.44	111.50
26	1H	2566	A	P-O3'-C3'	5.88	126.76	119.70
26	14	863	A	OP2-P-O3'	5.88	118.14	105.20
26	14	2281	C	N3-C4-N4	5.88	122.12	118.00
26	14	2287	A	N3-C4-C5	5.88	130.92	126.80
26	1H	1160	G	N7-C8-N9	5.88	116.04	113.10
26	1H	1357	U	N3-C4-C5	-5.88	111.07	114.60
26	1H	1319	G	C8-N9-C4	-5.88	104.05	106.40
26	1H	1902	C	C5-C4-N4	5.88	124.31	120.20
26	1H	2604	U	C6-N1-C2	-5.88	117.47	121.00
27	16	77	U	C5-C4-O4	-5.88	122.37	125.90
26	14	2392	A	O5'-P-OP1	-5.88	100.41	105.70
26	1H	581	C	N1-C2-O2	-5.88	115.37	118.90
26	1H	1915	U	N3-C4-C5	5.88	118.13	114.60
26	1H	26	G	N3-C4-N9	5.88	129.53	126.00
26	14	2719	G	C8-N9-C4	-5.88	104.05	106.40
26	14	791	C	N3-C2-O2	5.87	126.01	121.90
26	14	1616	A	C8-N9-C4	-5.87	103.45	105.80
1	13	857	C	C6-N1-C2	-5.87	117.95	120.30
26	1H	197	A	OP2-P-O3'	5.87	118.12	105.20
26	1H	564	C	C6-N1-C2	-5.87	117.95	120.30
26	1H	739	G	O5'-P-OP1	5.87	117.75	110.70
26	1H	795	C	O5'-P-OP2	-5.87	100.42	105.70
26	1H	1401	G	N7-C8-N9	5.87	116.04	113.10
26	1H	2727	G	N1-C6-O6	5.87	123.42	119.90
1	1G	769	G	N3-C4-C5	-5.87	125.66	128.60
26	14	1402	C	C6-N1-C2	-5.87	117.95	120.30
1	13	912	C	C6-N1-C2	5.87	122.65	120.30
26	14	2029	G	N1-C6-O6	5.87	123.42	119.90
26	14	2512	C	C4-C5-C6	-5.87	114.47	117.40
26	1H	2375	G	C5-C6-O6	-5.87	125.08	128.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	2530	A	N1-C6-N6	5.87	122.12	118.60
26	1H	2618	G	N9-C4-C5	5.87	107.75	105.40
26	14	2617	C	N3-C4-C5	5.87	124.25	121.90
26	1H	694	U	O5'-P-OP1	5.87	117.74	110.70
26	1H	1939	U	O5'-P-OP1	-5.87	100.42	105.70
26	1H	1200	C	C4-C5-C6	5.87	120.33	117.40
26	1H	1614	A	O4'-C1'-N9	5.87	112.89	108.20
26	1H	2508	G	N3-C2-N2	-5.87	115.79	119.90
1	1G	1259	C	C6-N1-C2	-5.87	117.95	120.30
26	14	527	C	C4-C5-C6	5.87	120.33	117.40
26	14	1320	C	N3-C4-N4	5.87	122.11	118.00
26	1H	845	G	C4-C5-N7	5.86	113.15	110.80
26	1H	970	C	C4-C5-C6	5.86	120.33	117.40
26	1H	1837	C	N1-C2-O2	5.86	122.42	118.90
26	1H	2578	G	N7-C8-N9	-5.86	110.17	113.10
1	1G	117	G	C5-C6-O6	-5.86	125.08	128.60
26	14	690	G	N1-C6-O6	5.86	123.42	119.90
26	14	1594	G	O5'-P-OP2	5.86	117.74	110.70
26	14	2526	G	N3-C4-C5	5.86	131.53	128.60
26	1H	1637	A	N1-C6-N6	-5.86	115.08	118.60
26	1H	1021	A	C8-N9-C4	-5.86	103.46	105.80
26	1H	2231	C	C6-N1-C2	-5.86	117.96	120.30
1	1G	17	U	C5-C6-N1	5.86	125.63	122.70
26	1H	777	A	C2-N3-C4	-5.86	107.67	110.60
26	1H	1828	G	C5-C6-O6	5.86	132.12	128.60
26	14	1394	U	C5-C6-N1	5.86	125.63	122.70
26	14	2699	C	N1-C2-O2	-5.86	115.39	118.90
1	13	760	G	C4-C5-N7	5.86	113.14	110.80
26	1H	1379	A	C2-N3-C4	-5.86	107.67	110.60
1	1G	1195	C	C6-N1-C2	-5.86	117.96	120.30
26	14	1926	U	N1-C2-N3	5.86	118.41	114.90
26	1H	818	G	C6-C5-N7	5.85	133.91	130.40
1	1G	690	G	C8-N9-C4	-5.85	104.06	106.40
26	1H	931	G	N1-C6-O6	-5.85	116.39	119.90
26	1H	954	G	OP2-P-O3'	5.85	118.08	105.20
26	1H	2243	U	O5'-P-OP2	-5.85	100.43	105.70
27	16	14	U	O5'-P-OP2	-5.85	100.43	105.70
1	1G	1119	C	C6-N1-C2	-5.85	117.96	120.30
26	14	2427	C	C6-N1-C2	5.85	122.64	120.30
26	1H	767	U	C5-C4-O4	5.85	129.41	125.90
26	1H	866	A	N7-C8-N9	5.85	116.72	113.80
24	3L	76	A	N1-C6-N6	5.85	122.11	118.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	1022	G	N9-C4-C5	5.85	107.74	105.40
26	14	1570	A	C6-C5-N7	-5.85	128.20	132.30
26	1H	1274	A	N7-C8-N9	5.85	116.72	113.80
1	1G	528	C	C5'-C4'-O4'	5.85	116.12	109.10
1	1G	800	G	N1-C6-O6	5.85	123.41	119.90
26	14	307	G	N1-C6-O6	5.85	123.41	119.90
26	14	2329	G	C5-C6-N1	5.85	114.42	111.50
26	1H	600	G	N1-C6-O6	5.85	123.41	119.90
27	16	49	C	N3-C4-N4	5.85	122.09	118.00
28	11	54	ARG	NE-CZ-NH2	-5.85	117.38	120.30
26	14	1804	C	O5'-P-OP2	5.85	117.72	110.70
1	1G	323	U	N3-C2-O2	5.85	126.29	122.20
26	14	1289	C	O5'-P-OP1	-5.85	100.44	105.70
1	13	276	G	N3-C2-N2	-5.84	115.81	119.90
1	13	771	G	C5-C6-O6	5.84	132.11	128.60
26	1H	936	C	C6-N1-C2	5.84	122.64	120.30
1	13	1495	U	O5'-P-OP2	-5.84	100.44	105.70
26	1H	1663	C	C6-N1-C2	5.84	122.64	120.30
26	1H	2297	C	O5'-P-OP2	-5.84	100.44	105.70
26	1H	2639	A	C2-N3-C4	-5.84	107.68	110.60
26	14	1989	G	N3-C2-N2	-5.84	115.81	119.90
26	14	2551	C	O5'-P-OP2	-5.84	100.44	105.70
26	14	2731	G	C4-N9-Cl'	5.84	134.10	126.50
26	1H	417	C	O5'-P-OP2	5.84	117.71	110.70
26	1H	729	G	C4-N9-Cl'	5.84	134.09	126.50
26	1H	816	C	OP1-P-OP2	-5.84	110.84	119.60
26	1H	1825	A	N9-C4-C5	5.84	108.14	105.80
26	1H	2259	G	O5'-P-OP2	5.84	117.71	110.70
26	1H	2310	A	C8-N9-C4	-5.84	103.46	105.80
27	16	23	G	N3-C4-C5	5.84	131.52	128.60
26	14	666	G	C2-N3-C4	-5.84	108.98	111.90
26	1H	113	G	OP1-P-O3'	5.84	118.05	105.20
26	1H	974(A)	C	P-O3'-C3'	5.84	126.71	119.70
26	1H	2594	C	C2-N3-C4	-5.84	116.98	119.90
26	14	1463	C	C6-N1-C2	-5.84	117.97	120.30
26	14	2571	C	N1-C2-O2	-5.84	115.40	118.90
26	14	2732	G	C5-C6-N1	5.84	114.42	111.50
26	14	1323	U	N3-C4-C5	-5.84	111.10	114.60
26	14	2017	U	O5'-P-OP1	-5.84	100.45	105.70
26	1H	103	A	C8-N9-C4	5.84	108.14	105.80
26	1H	1698	A	C6-C5-N7	-5.84	128.21	132.30
26	14	1422	G	N1-C6-O6	5.84	123.40	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	2441	C	N3-C4-N4	-5.84	113.92	118.00
26	1H	866	A	C5-N7-C8	-5.83	100.98	103.90
26	14	1446	C	C6-N1-C2	-5.83	117.97	120.30
1	13	14	U	O5'-P-OP1	-5.83	100.45	105.70
23	2K	6	G	N9-C4-C5	-5.83	103.07	105.40
26	1H	1678	G	N1-C2-N2	-5.83	110.95	116.20
26	1H	2429	G	OP1-P-OP2	-5.83	110.85	119.60
26	1H	2698	U	C4-C5-C6	5.83	123.20	119.70
26	14	71	A	C4-C5-N7	5.83	113.62	110.70
26	14	2387	U	C5-C6-N1	-5.83	119.78	122.70
26	1H	110	G	OP1-P-OP2	5.83	128.34	119.60
26	14	579	G	C5-C6-O6	-5.83	125.10	128.60
26	1H	792	G	N3-C4-N9	5.83	129.50	126.00
26	1H	1496	A	C5-C6-N6	-5.83	119.04	123.70
26	1H	2060	A	O4'-C1'-N9	5.83	112.86	108.20
26	14	2451	A	N7-C8-N9	5.83	116.71	113.80
26	14	2448	A	C5-C6-N1	5.83	120.61	117.70
26	14	119	A	OP1-P-O3'	5.82	118.01	105.20
26	14	1985	G	O5'-P-OP2	-5.82	100.46	105.70
1	13	963	G	N1-C6-O6	-5.82	116.41	119.90
1	13	610	G	N3-C4-N9	5.82	129.49	126.00
26	1H	1357	U	C5-C4-O4	5.82	129.39	125.90
26	1H	2688	U	C4-C5-C6	5.82	123.19	119.70
26	1H	1254	A	C4-C5-N7	5.82	113.61	110.70
26	1H	2303	G	OP1-P-O3'	5.82	118.00	105.20
26	14	409	C	C6-N1-C2	5.82	122.63	120.30
26	14	1643	G	OP2-P-O3'	5.82	118.00	105.20
26	1H	445	C	N3-C2-O2	-5.82	117.83	121.90
26	1H	733	G	N7-C8-N9	5.82	116.01	113.10
26	14	123	G	N1-C6-O6	5.82	123.39	119.90
26	14	2287	A	C5-C6-N1	-5.82	114.79	117.70
26	1H	2594	C	C4-C5-C6	5.81	120.31	117.40
26	14	198	C	O5'-P-OP1	-5.81	100.47	105.70
23	2K	1	C	C5-C6-N1	5.81	123.91	121.00
26	1H	318	C	O5'-P-OP1	-5.81	100.47	105.70
26	1H	397	G	C4-N9-C1'	-5.81	118.94	126.50
26	1H	818	G	N3-C4-N9	-5.81	122.51	126.00
26	14	2859	G	N9-C4-C5	5.81	107.72	105.40
25	4K	16	A	C8-N9-C4	5.81	108.12	105.80
1	13	509	A	P-O3'-C3'	5.81	126.67	119.70
26	1H	2228	G	N3-C4-C5	-5.81	125.69	128.60
26	1H	2269	A	N9-C4-C5	-5.81	103.48	105.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
27	16	25	A	OP1-P-O3'	5.81	117.98	105.20
26	14	2066	C	C5-C6-N1	5.81	123.91	121.00
1	13	1332	A	C8-N9-C4	-5.81	103.48	105.80
26	1H	1246	A	C4-C5-C6	5.81	119.90	117.00
26	14	2449	U	N3-C4-O4	5.81	123.47	119.40
26	1H	258	G	N3-C2-N2	5.80	123.96	119.90
26	1H	397	G	N3-C4-N9	-5.80	122.52	126.00
26	1H	660	G	C5-N7-C8	-5.80	101.40	104.30
1	1G	1354	C	C5-C6-N1	5.80	123.90	121.00
26	14	820	A	N1-C6-N6	-5.80	115.12	118.60
26	1H	330	A	N7-C8-N9	5.80	116.70	113.80
26	1H	530	G	N1-C2-N2	-5.80	110.98	116.20
1	1G	377	G	N3-C4-N9	5.80	129.48	126.00
26	1H	251	A	O5'-P-OP1	-5.80	100.48	105.70
26	14	2249	U	C5-C4-O4	5.80	129.38	125.90
26	14	2423	U	C5-C6-N1	-5.80	119.80	122.70
1	13	1508	G	N9-C4-C5	5.80	107.72	105.40
26	1H	2293	C	C6-N1-C2	-5.80	117.98	120.30
26	1H	2491	U	N3-C2-O2	5.80	126.26	122.20
26	14	97	C	OP1-P-OP2	5.80	128.30	119.60
26	14	630	G	C8-N9-C4	5.80	108.72	106.40
26	14	970	C	N1-C2-O2	-5.80	115.42	118.90
26	14	1391	U	C2-N1-C1'	5.80	124.66	117.70
26	14	2071	A	C8-N9-C4	-5.80	103.48	105.80
26	1H	785	G	C8-N9-C4	-5.80	104.08	106.40
26	1H	910	A	O5'-P-OP2	-5.80	100.48	105.70
26	1H	940	G	C5-C6-O6	-5.80	125.12	128.60
26	1H	1021	A	C5-C6-N1	-5.80	114.80	117.70
26	1H	1611	C	C4-C5-C6	5.80	120.30	117.40
26	14	2385	C	C5-C4-N4	-5.80	116.14	120.20
26	1H	488	G	O5'-P-OP2	-5.79	100.48	105.70
26	1H	1426	G	N1-C6-O6	-5.79	116.42	119.90
26	14	778	G	N1-C6-O6	-5.79	116.42	119.90
23	2K	17	C	C2-N1-C1'	5.79	125.17	118.80
26	1H	1496	A	O4'-C1'-N9	5.79	112.83	108.20
26	1H	1761	C	N1-C2-O2	-5.79	115.42	118.90
26	1H	1773	A	C5-C6-N1	-5.79	114.80	117.70
26	1H	1932	A	N1-C6-N6	5.79	122.08	118.60
1	1G	1502	A	N1-C2-N3	5.79	132.20	129.30
26	14	1029	A	N9-C4-C5	-5.79	103.48	105.80
1	13	1219	U	C6-N1-C2	-5.79	117.53	121.00
26	1H	1336	A	C5-C6-N1	5.79	120.60	117.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	13	680	C	O5'-P-OP2	-5.79	100.49	105.70
26	14	818	G	N7-C8-N9	5.79	116.00	113.10
26	14	1396	U	C2-N1-C1'	5.79	124.65	117.70
1	13	580	U	C6-N1-C2	5.79	124.47	121.00
26	1H	1525	G	O5'-P-OP2	-5.79	100.49	105.70
26	14	1216	G	C4-N9-C1'	5.79	134.02	126.50
26	14	1519	G	C4-C5-N7	-5.79	108.48	110.80
26	1H	676	A	OP1-P-OP2	5.79	128.28	119.60
26	1H	1399	C	OP2-P-O3'	5.79	117.93	105.20
26	14	833	U	N1-C2-O2	-5.79	118.75	122.80
26	14	1470	G	N1-C6-O6	5.79	123.37	119.90
1	13	1502	A	N9-C4-C5	-5.79	103.49	105.80
26	1H	1828	G	C6-N1-C2	5.79	128.57	125.10
26	14	530	G	C8-N9-C1'	-5.79	119.48	127.00
26	14	783	A	O5'-P-OP1	5.79	117.64	110.70
26	1H	1013	C	N3-C2-O2	5.78	125.95	121.90
26	14	388	G	C2-N3-C4	-5.78	109.01	111.90
26	14	678	C	N3-C4-C5	5.78	124.21	121.90
26	14	1328	G	N9-C4-C5	-5.78	103.09	105.40
26	14	1379	A	C5-C6-N6	-5.78	119.07	123.70
1	13	756	C	C4-C5-C6	5.78	120.29	117.40
26	1H	1011	G	O5'-P-OP1	-5.78	100.50	105.70
26	1H	1246	A	C6-N1-C2	-5.78	115.13	118.60
26	1H	1851	U	N1-C2-O2	-5.78	118.75	122.80
26	1H	2492	U	N3-C2-O2	-5.78	118.15	122.20
26	14	2240	C	C5-C6-N1	5.78	123.89	121.00
26	1H	928	G	N3-C4-C5	5.78	131.49	128.60
26	1H	1938	A	O5'-P-OP1	-5.78	100.50	105.70
26	1H	2004	G	C6-C5-N7	-5.78	126.93	130.40
26	1H	2060	A	C4-C5-C6	-5.78	114.11	117.00
4	32	157	LEU	CA-CB-CG	5.78	128.59	115.30
26	14	1225	C	C6-N1-C2	5.78	122.61	120.30
26	14	1323	U	OP1-P-OP2	-5.78	110.93	119.60
26	1H	2070	G	C5-C6-O6	5.78	132.07	128.60
1	13	731	G	OP1-P-O3'	5.78	117.91	105.20
1	13	1469	G	N7-C8-N9	5.78	115.99	113.10
26	1H	676	A	C6-N1-C2	5.78	122.07	118.60
26	1H	940	G	C6-C5-N7	-5.78	126.93	130.40
26	14	265	A	N1-C6-N6	5.78	122.07	118.60
26	14	1189	A	OP1-P-OP2	-5.78	110.93	119.60
1	13	1359	C	O5'-P-OP1	-5.78	100.50	105.70
26	1H	1927	A	C8-N9-C4	-5.78	103.49	105.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	2554	U	C5-C4-O4	-5.78	122.43	125.90
26	14	201	C	C6-N1-C2	5.78	122.61	120.30
26	14	2573	C	C5-C6-N1	5.78	123.89	121.00
26	14	1468	C	C5-C6-N1	5.77	123.89	121.00
26	14	2392	A	N7-C8-N9	5.77	116.69	113.80
1	13	813	U	OP1-P-OP2	5.77	128.26	119.60
26	1H	906	G	N1-C6-O6	-5.77	116.44	119.90
26	1H	1298	C	C2-N3-C4	5.77	122.79	119.90
26	1H	1347	G	C5-C6-O6	-5.77	125.14	128.60
26	1H	199	A	C4-C5-C6	-5.77	114.11	117.00
1	1G	121	C	C2-N1-C1'	5.77	125.15	118.80
26	14	1396	U	N1-C2-O2	5.77	126.84	122.80
26	14	1519	G	C5-C6-O6	5.77	132.06	128.60
26	14	2429	G	OP2-P-O3'	5.77	117.89	105.20
26	1H	961	C	O4'-C1'-N1	5.77	112.81	108.20
26	1H	1217	C	O5'-P-OP1	-5.77	100.51	105.70
26	1H	1626	G	C6-C5-N7	-5.77	126.94	130.40
26	1H	2068	U	N1-C2-O2	5.77	126.84	122.80
54	P8	9	ARG	NE-CZ-NH2	5.77	123.18	120.30
1	1G	137	C	C6-N1-C2	5.77	122.61	120.30
26	1H	1776	G	C4-C5-N7	5.77	113.11	110.80
26	14	577	G	C5-C6-O6	-5.77	125.14	128.60
26	14	2251	G	N3-C4-C5	-5.77	125.72	128.60
26	1H	1626	G	C5-N7-C8	-5.76	101.42	104.30
26	14	2606	C	OP2-P-O3'	5.76	117.88	105.20
26	1H	461	C	N3-C2-O2	5.76	125.93	121.90
26	1H	1950	G	N3-C2-N2	5.76	123.93	119.90
26	1H	2051	A	C8-N9-C4	-5.76	103.50	105.80
1	1G	748	C	P-O3'-C3'	5.76	126.62	119.70
26	14	778	G	C5-C6-O6	5.76	132.06	128.60
26	1H	752	A	N1-C2-N3	5.76	132.18	129.30
26	14	969	U	C6-N1-C2	-5.76	117.54	121.00
26	14	1965	C	C2-N1-C1'	-5.76	112.46	118.80
1	13	745	C	C5-C6-N1	5.76	123.88	121.00
1	13	963	G	N3-C2-N2	5.76	123.93	119.90
26	1H	461	C	C4-C5-C6	5.76	120.28	117.40
26	1H	793	A	N3-C4-N9	5.76	132.01	127.40
26	1H	2584	U	N1-C2-O2	5.76	126.83	122.80
26	14	697	C	C6-N1-C2	-5.76	118.00	120.30
26	14	1333	C	N3-C4-C5	5.76	124.20	121.90
26	1H	647	G	C8-N9-C4	-5.76	104.10	106.40
26	14	123	G	OP1-P-OP2	5.76	128.24	119.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	13	529	G	N9-C4-C5	-5.76	103.10	105.40
1	13	623	C	C5-C6-N1	5.76	123.88	121.00
1	1G	501	C	C6-N1-C2	-5.76	118.00	120.30
1	1G	1446	A	O4'-C1'-N9	5.76	112.81	108.20
1	13	373	A	C8-N9-C4	5.75	108.10	105.80
26	1H	192	C	C5-C4-N4	-5.75	116.17	120.20
26	1H	1950	G	C6-C5-N7	-5.75	126.95	130.40
1	13	891	U	OP2-P-O3'	5.75	117.86	105.20
1	13	1475	G	C8-N9-C4	-5.75	104.10	106.40
26	1H	116	C	N1-C2-O2	-5.75	115.45	118.90
26	1H	1224	G	N3-C4-C5	5.75	131.48	128.60
26	1H	1415	U	C5-C4-O4	5.75	129.35	125.90
26	1H	1544	C	N1-C2-O2	5.75	122.35	118.90
26	14	2062	A	C8-N9-C4	5.75	108.10	105.80
26	14	2678	C	C6-N1-C2	5.75	122.60	120.30
1	13	826	C	C6-N1-C2	-5.75	118.00	120.30
1	13	872	A	N1-C2-N3	-5.75	126.42	129.30
26	1H	119	A	C4-C5-C6	5.75	119.88	117.00
26	1H	865	C	C6-N1-C2	5.75	122.60	120.30
26	1H	2766	G	C5-C6-O6	-5.75	125.15	128.60
1	1G	197	A	C8-N9-C4	-5.75	103.50	105.80
26	14	1771	C	C5-C4-N4	-5.75	116.17	120.20
26	14	1774	C	O5'-P-OP2	5.75	117.60	110.70
26	14	1962	C	C5-C4-N4	-5.75	116.17	120.20
26	14	2685	G	C8-N9-C4	5.75	108.70	106.40
26	1H	139	G	N3-C4-C5	-5.75	125.72	128.60
26	1H	203	C	N3-C2-O2	5.75	125.92	121.90
26	1H	1284	A	C2-N3-C4	-5.75	107.73	110.60
26	1H	1287	A	O5'-P-OP1	5.75	117.60	110.70
26	1H	2392	A	C6-C5-N7	-5.75	128.28	132.30
23	2L	21	U	C4-C5-C6	-5.75	116.25	119.70
1	13	1501	C	C6-N1-C2	-5.75	118.00	120.30
26	1H	198	C	N3-C2-O2	-5.75	117.88	121.90
26	1H	768	G	C5-C6-O6	5.75	132.05	128.60
26	1H	1626	G	N3-C2-N2	-5.75	115.88	119.90
26	14	130	C	C2-N3-C4	-5.75	117.03	119.90
26	14	528	A	N1-C2-N3	5.75	132.17	129.30
26	1H	1312	U	C5-C4-O4	5.75	129.35	125.90
26	1H	2396	G	C4-N9-C1'	-5.75	119.03	126.50
26	14	1605	C	N1-C2-O2	-5.75	115.45	118.90
26	1H	128	C	N3-C4-C5	5.74	124.20	121.90
26	1H	954	G	C4-C5-N7	-5.74	108.50	110.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	1347	G	N1-C6-O6	5.74	123.35	119.90
27	1J	22	U	C5-C6-N1	5.74	125.57	122.70
26	1H	196	A	C6-N1-C2	5.74	122.05	118.60
26	1H	2007	C	O5'-P-OP2	-5.74	100.53	105.70
26	1H	2287	A	N3-C4-C5	5.74	130.82	126.80
26	14	752	A	N1-C2-N3	5.74	132.17	129.30
1	13	644	G	C8-N9-C4	5.74	108.70	106.40
26	1H	802	A	C5-N7-C8	-5.74	101.03	103.90
26	1H	1764	G	N1-C6-O6	-5.74	116.46	119.90
23	2L	50	G	C8-N9-C4	-5.74	104.10	106.40
26	14	252	G	N9-C4-C5	5.74	107.70	105.40
26	14	270(Y)	G	C5-C6-O6	5.74	132.04	128.60
26	14	827	U	N1-C2-O2	-5.74	118.78	122.80
26	14	2432	A	N1-C6-N6	5.74	122.05	118.60
1	13	1426	C	N3-C4-N4	5.74	122.02	118.00
26	1H	533	G	N1-C2-N3	5.74	127.34	123.90
26	1H	845	G	C5-N7-C8	-5.74	101.43	104.30
26	1H	1814	G	C8-N9-C4	5.74	108.69	106.40
26	1H	2497	A	N1-C2-N3	5.74	132.17	129.30
26	1H	2592	G	C5-C6-N1	-5.74	108.63	111.50
26	14	2607	G	O5'-P-OP1	5.74	117.59	110.70
36	35	62	LEU	CA-CB-CG	-5.74	102.10	115.30
26	1H	1299	G	C4-C5-N7	5.74	113.09	110.80
26	14	191	A	N1-C6-N6	5.74	122.04	118.60
26	1H	209	C	N3-C4-C5	5.74	124.19	121.90
26	14	1312	U	O5'-P-OP1	-5.74	100.54	105.70
26	1H	471	A	C5-N7-C8	-5.73	101.03	103.90
26	1H	451	C	C6-N1-C2	5.73	122.59	120.30
26	1H	735	A	C6-N1-C2	-5.73	115.16	118.60
26	1H	1622	G	N9-C4-C5	5.73	107.69	105.40
26	1H	1622	G	C6-C5-N7	5.73	133.84	130.40
26	1H	2018	G	N7-C8-N9	5.73	115.97	113.10
26	1H	2718	G	C5-C6-O6	-5.73	125.16	128.60
26	14	1600	C	C6-N1-C2	5.73	122.59	120.30
26	14	2501	C	C6-N1-C2	5.73	122.59	120.30
1	13	365	U	C2-N1-C1'	5.73	124.58	117.70
26	1H	2446	G	C6-C5-N7	-5.73	126.96	130.40
26	14	205	G	N3-C4-N9	5.73	129.44	126.00
26	14	252	G	C6-C5-N7	5.73	133.84	130.40
26	14	855	G	N7-C8-N9	5.73	115.97	113.10
26	14	1314	C	N1-C2-O2	5.73	122.34	118.90
1	13	749	C	C2-N1-C1'	5.73	125.10	118.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	2702	U	N1-C1'-C2'	5.73	121.45	114.00
26	1H	779	U	C6-N1-C2	5.73	124.44	121.00
26	14	2401	U	N3-C4-O4	5.73	123.41	119.40
26	14	2763	G	N1-C6-O6	-5.73	116.46	119.90
26	1H	141	A	C4-C5-N7	5.73	113.56	110.70
26	1H	792	G	N1-C6-O6	5.73	123.34	119.90
26	1H	972	G	C6-C5-N7	5.73	133.84	130.40
26	1H	2079	U	N1-C2-O2	-5.73	118.79	122.80
26	14	2433	A	N7-C8-N9	5.73	116.66	113.80
24	3K	71	G	C4-C5-N7	-5.72	108.51	110.80
26	1H	300	A	O4'-C1'-N9	5.72	112.78	108.20
26	1H	972	G	O5'-P-OP1	5.72	117.57	110.70
26	1H	2827	C	N3-C2-O2	5.72	125.91	121.90
26	14	694	U	O5'-P-OP2	-5.72	100.55	105.70
26	14	1827	C	N1-C2-O2	5.72	122.33	118.90
26	14	1905	C	O5'-P-OP2	-5.72	100.55	105.70
26	1H	708	C	O5'-P-OP1	5.72	117.57	110.70
26	1H	744	G	O5'-P-OP2	-5.72	100.55	105.70
26	1H	2210	G	C4-N9-C1'	5.72	133.94	126.50
1	1G	789	U	C6-N1-C2	-5.72	117.57	121.00
26	14	74	A	N1-C2-N3	5.72	132.16	129.30
26	14	861	A	O5'-P-OP2	5.72	117.57	110.70
26	14	2873	A	C5-C6-N1	-5.72	114.84	117.70
26	1H	996	A	N7-C8-N9	-5.72	110.94	113.80
26	14	684	G	N9-C4-C5	5.72	107.69	105.40
26	14	1992	G	C8-N9-C4	5.72	108.69	106.40
26	14	2818	G	N1-C6-O6	5.72	123.33	119.90
26	1H	845	G	C4-N9-C1'	-5.72	119.06	126.50
26	1H	866	A	C6-C5-N7	-5.72	128.30	132.30
26	1H	1936	A	N1-C6-N6	5.72	122.03	118.60
26	1H	2310	A	N7-C8-N9	5.72	116.66	113.80
26	1H	2702	U	C5'-C4'-O4'	5.72	115.96	109.10
1	1G	1432	G	N7-C8-N9	5.72	115.96	113.10
26	14	476	G	OP1-P-OP2	5.72	128.18	119.60
26	14	1386	C	C6-N1-C2	-5.72	118.01	120.30
26	14	1982	C	C2-N1-C1'	5.72	125.09	118.80
26	1H	1653	G	N1-C2-N2	-5.72	111.05	116.20
26	1H	2466	C	C6-N1-C2	5.72	122.59	120.30
26	1H	2581	G	N3-C2-N2	5.72	123.90	119.90
26	14	2492	U	O5'-P-OP2	5.72	117.56	110.70
1	13	266	G	C5-N7-C8	-5.72	101.44	104.30
1	13	1468	A	N7-C8-N9	-5.72	110.94	113.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	464	U	O5'-P-OP1	-5.72	100.56	105.70
26	1H	533	G	N3-C4-C5	-5.72	125.74	128.60
26	1H	1299	G	C5-N7-C8	-5.72	101.44	104.30
26	1H	1400	G	C8-N9-C4	-5.72	104.11	106.40
26	14	106	C	C6-N1-C2	-5.72	118.01	120.30
26	14	1528	A	N7-C8-N9	5.72	116.66	113.80
26	14	2251	G	C4-C5-N7	-5.72	108.51	110.80
26	14	2873	A	O5'-P-OP1	-5.72	100.56	105.70
1	13	1498	U	C5-C4-O4	-5.71	122.47	125.90
26	1H	2082	A	C8-N9-C4	5.71	108.09	105.80
26	14	667	U	OP2-P-O3'	5.71	117.77	105.20
26	14	2365	G	N3-C4-N9	5.71	129.43	126.00
26	14	2525	G	OP2-P-O3'	5.71	117.77	105.20
1	13	1446	A	O4'-C1'-N9	5.71	112.77	108.20
26	14	1742	C	C5-C6-N1	5.71	123.86	121.00
26	1H	1107	G	C8-N9-C4	-5.71	104.12	106.40
26	1H	1332	G	O4'-C1'-N9	-5.71	103.63	108.20
26	1H	1470	G	OP2-P-O3'	5.71	117.76	105.20
26	14	1681	G	N3-C4-C5	5.71	131.45	128.60
26	1H	1769	G	O5'-P-OP2	-5.71	100.56	105.70
26	1H	2490	G	C8-N9-C4	-5.71	104.12	106.40
26	1H	2700	C	C5-C6-N1	-5.71	118.15	121.00
1	1G	1270	C	C6-N1-C2	-5.71	118.02	120.30
26	14	577	G	O5'-P-OP2	5.71	117.55	110.70
26	14	929	G	C4-C5-C6	5.71	122.22	118.80
26	14	1029	A	C5-C6-N6	-5.71	119.14	123.70
26	14	1980	G	N1-C6-O6	5.71	123.32	119.90
26	14	2702	U	C2-N1-C1'	5.71	124.55	117.70
26	1H	71	A	N3-C4-C5	5.70	130.79	126.80
26	1H	948	G	O5'-P-OP2	5.70	117.54	110.70
26	1H	1013	C	N1-C2-O2	-5.70	115.48	118.90
26	1H	1624	G	C8-N9-C4	5.70	108.68	106.40
26	14	562	U	N3-C2-O2	-5.70	118.21	122.20
26	14	1255	U	N3-C2-O2	-5.70	118.21	122.20
26	14	2063	C	OP2-P-O3'	5.70	117.75	105.20
26	1H	2346	A	N1-C6-N6	5.70	122.02	118.60
1	13	332	G	O5'-P-OP1	-5.70	100.57	105.70
22	1K	18	G	C8-N9-C1'	-5.70	119.59	127.00
26	1H	1518	C	OP1-P-O3'	5.70	117.74	105.20
26	14	2866	U	O5'-P-OP2	-5.70	100.57	105.70
26	1H	192	C	C5-C6-N1	5.70	123.85	121.00
26	1H	498	G	O5'-P-OP2	5.70	117.54	110.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	1G	322	C	C5-C4-N4	-5.70	116.21	120.20
1	1G	900	A	N1-C6-N6	5.70	122.02	118.60
26	14	530	G	N3-C4-C5	5.70	131.45	128.60
26	14	1642	G	OP2-P-O3'	5.70	117.74	105.20
26	1H	28	A	N7-C8-N9	-5.70	110.95	113.80
26	1H	793	A	C6-C5-N7	-5.70	128.31	132.30
26	1H	2060	A	C6-C5-N7	5.70	136.29	132.30
23	2K	48	U	P-O3'-C3'	5.70	126.53	119.70
26	1H	557	U	C5-C6-N1	-5.70	119.85	122.70
26	1H	755	C	C4-C5-C6	5.70	120.25	117.40
26	1H	2318	G	C5-N7-C8	-5.70	101.45	104.30
1	1G	1527	C	C2-N3-C4	-5.70	117.05	119.90
26	14	733	G	N7-C8-N9	5.70	115.95	113.10
26	14	2440	C	N1-C2-N3	-5.70	115.21	119.20
26	14	512	G	C5-C6-O6	5.69	132.02	128.60
1	13	1112	C	C6-N1-C2	-5.69	118.02	120.30
26	1H	627	A	C8-N9-C4	5.69	108.08	105.80
26	1H	1902	C	N3-C4-N4	-5.69	114.02	118.00
26	14	602	G	N3-C4-N9	5.69	129.42	126.00
1	13	584	G	N1-C6-O6	-5.69	116.48	119.90
26	1H	470	A	C5-N7-C8	-5.69	101.06	103.90
26	1H	1936	A	N7-C8-N9	5.69	116.65	113.80
26	1H	2323	G	C8-N9-C4	-5.69	104.12	106.40
26	1H	203	C	C2-N1-C1'	-5.69	112.54	118.80
26	1H	224	G	O5'-P-OP2	-5.69	100.58	105.70
26	1H	1779	U	OP1-P-O3'	5.69	117.71	105.20
26	1H	2377	A	C2-N3-C4	-5.69	107.76	110.60
26	1H	792	G	N3-C4-C5	-5.69	125.76	128.60
26	1H	2406	U	N3-C2-O2	-5.69	118.22	122.20
27	16	77	U	C2-N3-C4	-5.69	123.59	127.00
26	1H	558	G	O5'-P-OP2	-5.68	100.58	105.70
23	2L	50	G	N3-C4-C5	-5.68	125.76	128.60
26	14	2449	U	C5-C4-O4	-5.68	122.49	125.90
26	1H	1241	A	C2-N3-C4	-5.68	107.76	110.60
26	1H	1305	C	N3-C2-O2	-5.68	117.92	121.90
26	1H	2447	G	N3-C4-C5	-5.68	125.76	128.60
1	1G	1205	U	C5-C6-N1	5.68	125.54	122.70
26	14	1694	C	N1-C2-O2	5.68	122.31	118.90
26	1H	457	A	N1-C6-N6	-5.68	115.19	118.60
26	14	2607	G	C6-C5-N7	-5.68	126.99	130.40
26	1H	470	A	C2-N3-C4	-5.68	107.76	110.60
26	1H	771	G	C5-C6-O6	-5.68	125.19	128.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	793	A	C4-C5-C6	5.68	119.84	117.00
26	1H	839	U	N1-C2-O2	-5.68	118.82	122.80
26	1H	2377	A	N9-C4-C5	-5.68	103.53	105.80
26	1H	2428	G	N9-C4-C5	5.68	107.67	105.40
26	1H	2452	C	N3-C4-C5	5.68	124.17	121.90
27	16	44	G	C8-N9-C1'	5.68	134.38	127.00
1	1G	1048	G	C8-N9-C4	-5.68	104.13	106.40
26	14	1614	A	O4'-C1'-N9	5.68	112.74	108.20
26	14	2459	A	N7-C8-N9	5.68	116.64	113.80
26	14	2730	C	C6-N1-C2	-5.68	118.03	120.30
26	1H	448	U	C5-C6-N1	-5.68	119.86	122.70
1	1G	1139	G	N3-C4-C5	5.68	131.44	128.60
26	14	1309	G	C8-N9-C1'	-5.68	119.62	127.00
26	1H	1370	C	C2-N1-C1'	-5.68	112.56	118.80
26	1H	1603	A	OP1-P-OP2	-5.68	111.09	119.60
26	1H	2300	G	N7-C8-N9	5.68	115.94	113.10
24	1L	74	C	C5-C6-N1	-5.68	118.16	121.00
26	14	1607	C	C5-C4-N4	-5.68	116.23	120.20
1	13	858	G	N3-C4-C5	-5.67	125.76	128.60
26	14	1367	A	N9-C4-C5	-5.67	103.53	105.80
26	1H	2433	A	O5'-P-OP2	5.67	117.51	110.70
26	14	93	C	C2-N1-C1'	5.67	125.04	118.80
26	14	1251	C	OP1-P-OP2	5.67	128.11	119.60
26	14	1283	G	OP1-P-OP2	5.67	128.11	119.60
26	14	1474	C	C5-C6-N1	5.67	123.84	121.00
26	1H	52	A	C2-N3-C4	5.67	113.44	110.60
26	1H	755	C	OP2-P-O3'	5.67	117.67	105.20
26	1H	1610	A	O4'-C1'-N9	-5.67	103.66	108.20
26	1H	1800	C	C6-N1-C2	-5.67	118.03	120.30
26	1H	1982	C	N3-C4-N4	5.67	121.97	118.00
26	1H	2871	C	O5'-P-OP2	-5.67	100.60	105.70
27	16	102	G	N3-C2-N2	-5.67	115.93	119.90
26	1H	1914	C	N1-C2-O2	5.67	122.30	118.90
26	1H	1940	U	C4-C5-C6	5.67	123.10	119.70
1	1G	1203	C	C6-N1-C2	5.67	122.57	120.30
26	14	1629	U	N3-C4-O4	5.67	123.37	119.40
26	14	1772	G	N1-C6-O6	5.67	123.30	119.90
1	13	858	G	C8-N9-C4	-5.67	104.13	106.40
26	1H	185	U	C5-C6-N1	-5.67	119.87	122.70
26	1H	211	A	C2-N3-C4	-5.67	107.77	110.60
26	1H	456	C	C5-C6-N1	-5.67	118.17	121.00
26	1H	825	C	N3-C4-N4	5.67	121.97	118.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	986	C	OP1-P-OP2	-5.67	111.10	119.60
26	1H	1258	C	O5'-P-OP2	-5.67	100.60	105.70
26	1H	1366	A	C8-N9-C4	5.67	108.07	105.80
26	1H	1885	A	N7-C8-N9	-5.67	110.97	113.80
26	1H	1913	A	N1-C6-N6	5.67	122.00	118.60
26	1H	47	C	C5-C4-N4	-5.67	116.23	120.20
26	1H	1970	A	N7-C8-N9	5.66	116.63	113.80
26	1H	2444	G	OP1-P-O3'	5.66	117.66	105.20
1	1G	328	C	P-O3'-C3'	5.66	126.50	119.70
1	13	1422	G	C8-N9-C4	5.66	108.67	106.40
26	1H	452	G	N1-C6-O6	-5.66	116.50	119.90
26	1H	1942	C	N3-C4-C5	5.66	124.16	121.90
26	14	746	A	O5'-P-OP1	-5.66	100.61	105.70
26	14	2712	U	C4-C5-C6	5.66	123.10	119.70
26	1H	1221	C	N3-C2-O2	-5.66	117.94	121.90
26	14	866	A	C8-N9-C1'	-5.66	117.52	127.70
26	14	982	C	C6-N1-C2	-5.66	118.04	120.30
26	1H	214	G	O5'-P-OP2	-5.66	100.61	105.70
27	16	24	G	C8-N9-C4	-5.66	104.14	106.40
26	1H	487	C	O5'-P-OP1	-5.66	100.61	105.70
26	1H	1944	U	N3-C2-O2	-5.66	118.24	122.20
1	1G	117	G	C6-C5-N7	-5.66	127.01	130.40
1	1G	442	C	C6-N1-C2	-5.66	118.04	120.30
26	1H	842	G	N3-C4-C5	5.65	131.43	128.60
26	1H	2710	C	OP2-P-O3'	5.65	117.64	105.20
26	14	1441	G	C8-N9-C4	5.65	108.66	106.40
26	1H	1310	G	C5-C6-O6	-5.65	125.21	128.60
26	1H	2036	C	OP2-P-O3'	5.65	117.63	105.20
26	1H	2503	A	OP2-P-O3'	5.65	117.64	105.20
26	14	1609	A	C6-N1-C2	-5.65	115.21	118.60
26	1H	67	U	C5-C6-N1	5.65	125.53	122.70
1	1G	894	G	C5-C6-O6	-5.65	125.21	128.60
26	14	574	C	N3-C4-N4	-5.65	114.04	118.00
26	1H	1782	C	C4-C5-C6	5.65	120.22	117.40
26	1H	2282	G	O5'-P-OP2	5.65	117.48	110.70
27	1J	113	C	O4'-C1'-N1	5.65	112.72	108.20
1	13	656	C	C2-N3-C4	5.65	122.72	119.90
26	1H	251	A	O5'-P-OP2	5.65	117.48	110.70
26	1H	1409	C	O5'-P-OP1	-5.65	100.62	105.70
26	14	114	U	C2-N1-C1'	5.65	124.48	117.70
1	13	509	A	C2'-C3'-O3'	5.65	122.73	113.70
26	1H	123	G	C5-C6-N1	5.65	114.32	111.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	300	A	O5'-P-OP2	-5.65	100.62	105.70
26	1H	688	U	OP1-P-OP2	5.65	128.07	119.60
26	1H	1558	A	P-O3'-C3'	5.65	126.47	119.70
26	14	747	U	OP1-P-OP2	5.65	128.07	119.60
26	14	1494	A	C8-N9-C4	-5.65	103.54	105.80
1	13	1245	A	O5'-P-OP2	-5.64	100.62	105.70
26	1H	572	A	C8-N9-C4	-5.64	103.54	105.80
26	1H	1309	G	O5'-P-OP1	5.64	117.47	110.70
26	1H	1837	C	O5'-P-OP1	-5.64	100.62	105.70
26	1H	2030	A	C5-C6-N6	-5.64	119.18	123.70
26	1H	127	A	C5-C6-N6	-5.64	119.19	123.70
26	1H	684	G	C8-N9-C4	-5.64	104.14	106.40
26	14	782	A	C6-N1-C2	-5.64	115.21	118.60
26	14	866	A	O4'-C1'-N9	-5.64	103.69	108.20
26	14	1266	G	C5-C6-O6	-5.64	125.21	128.60
26	14	1775	U	OP1-P-OP2	-5.64	111.14	119.60
26	14	840	C	N3-C2-O2	5.64	125.85	121.90
26	14	1325	G	C4-C5-N7	5.64	113.06	110.80
1	13	900	A	OP1-P-OP2	-5.64	111.14	119.60
26	1H	699	A	C2-N3-C4	5.64	113.42	110.60
26	14	690	G	OP1-P-O3'	5.64	117.61	105.20
26	14	2674	G	O5'-P-OP2	-5.64	100.62	105.70
1	13	960	U	C2-N1-C1'	5.64	124.47	117.70
26	1H	1372	U	OP2-P-O3'	5.64	117.60	105.20
26	14	2367	G	C8-N9-C4	-5.64	104.14	106.40
26	1H	667	U	N3-C2-O2	5.64	126.15	122.20
23	2L	35	C	C6-N1-C1'	-5.64	114.04	120.80
26	14	642	G	N7-C8-N9	5.64	115.92	113.10
26	14	2068	U	OP1-P-O3'	5.64	117.60	105.20
26	14	2072	G	C4-C5-N7	5.64	113.05	110.80
1	13	1409	C	C5-C4-N4	-5.63	116.26	120.20
26	1H	467	G	C8-N9-C4	5.63	108.65	106.40
26	1H	805	G	N3-C4-N9	5.63	129.38	126.00
27	16	7	G	C4-C5-N7	5.63	113.05	110.80
26	14	1768	U	C5-C4-O4	5.63	129.28	125.90
26	14	2351	G	C4-N9-C1'	5.63	133.83	126.50
26	1H	2070	G	N9-C4-C5	-5.63	103.15	105.40
1	13	1069	C	C6-N1-C2	-5.63	118.05	120.30
26	1H	1544	C	C6-N1-C2	5.63	122.55	120.30
26	14	562	U	N1-C2-N3	5.63	118.28	114.90
26	14	1026	U	C5-C6-N1	5.63	125.52	122.70
26	14	1516	U	N1-C2-O2	5.63	126.74	122.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	2826	A	N9-C4-C5	5.63	108.05	105.80
1	13	730	G	OP1-P-O3'	5.63	117.59	105.20
1	13	1246	C	C6-N1-C2	-5.63	118.05	120.30
26	1H	66	C	OP1-P-OP2	-5.63	111.15	119.60
26	1H	165	U	C2-N1-C1'	5.63	124.46	117.70
26	1H	756	C	N1-C2-O2	-5.63	115.52	118.90
26	1H	733	G	N3-C2-N2	5.63	123.84	119.90
26	14	1970	A	C8-N9-C4	-5.63	103.55	105.80
26	14	2039	C	C5-C6-N1	5.63	123.81	121.00
26	14	2333	A	O5'-P-OP2	-5.63	100.63	105.70
26	1H	1445	C	C2-N3-C4	5.63	122.71	119.90
26	14	1545(A)	A	C8-N9-C4	-5.63	103.55	105.80
26	14	2062	A	C5-C6-N6	-5.62	119.20	123.70
1	13	277	C	C6-N1-C2	-5.62	118.05	120.30
26	1H	1011	G	N3-C4-C5	-5.62	125.79	128.60
26	1H	1022	G	C8-N9-C4	-5.62	104.15	106.40
26	1H	1381	G	N3-C2-N2	-5.62	115.96	119.90
26	1H	2247	A	C2-N3-C4	-5.62	107.79	110.60
26	14	752	A	N7-C8-N9	5.62	116.61	113.80
26	14	818	G	N9-C4-C5	5.62	107.65	105.40
26	14	951	C	C6-N1-C2	-5.62	118.05	120.30
26	1H	1606	G	O5'-P-OP2	-5.62	100.64	105.70
26	1H	2280	G	C2-N3-C4	5.62	114.71	111.90
26	14	819	A	C4-C5-C6	5.62	119.81	117.00
26	1H	1593	G	OP1-P-O3'	5.62	117.56	105.20
26	14	700	G	OP1-P-OP2	-5.62	111.17	119.60
26	14	2592	G	N3-C4-N9	5.62	129.37	126.00
26	1H	1602	U	O5'-P-OP2	5.62	117.44	110.70
26	1H	2506	U	N3-C2-O2	-5.62	118.27	122.20
24	3L	76	A	C6-C5-N7	-5.62	128.37	132.30
26	14	1142	U	N1-C2-O2	5.62	126.73	122.80
1	13	326	G	C5-C6-O6	5.62	131.97	128.60
1	13	1359	C	C6-N1-C2	5.62	122.55	120.30
1	13	127	G	OP2-P-O3'	5.62	117.55	105.20
26	1H	183	C	N1-C2-O2	5.62	122.27	118.90
26	1H	430	G	N1-C6-O6	5.62	123.27	119.90
26	1H	746	A	N1-C6-N6	5.62	121.97	118.60
26	1H	1962	C	C2-N1-C1'	-5.62	112.62	118.80
26	1H	2689	U	C2-N3-C4	-5.62	123.63	127.00
26	14	1248	G	C8-N9-C4	5.62	108.65	106.40
26	14	1325	G	N1-C6-O6	5.62	123.27	119.90
1	13	767	A	C2-N3-C4	-5.61	107.79	110.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
23	2K	27	G	O5'-P-OP2	5.61	117.44	110.70
26	1H	566	U	N3-C4-C5	5.61	117.97	114.60
26	1H	1313	U	C2-N1-C1'	5.61	124.44	117.70
26	1H	2250	G	N9-C4-C5	5.61	107.65	105.40
27	16	81	G	C6-C5-N7	-5.61	127.03	130.40
26	14	791	C	C2-N3-C4	-5.61	117.09	119.90
26	1H	99	U	N3-C2-O2	-5.61	118.27	122.20
26	1H	183	C	N3-C2-O2	-5.61	117.97	121.90
26	1H	397	G	C8-N9-C1'	5.61	134.29	127.00
26	1H	2614	A	C2-N3-C4	5.61	113.41	110.60
26	14	821	A	OP1-P-OP2	5.61	128.02	119.60
26	14	2224	G	O5'-P-OP1	-5.61	100.65	105.70
26	1H	16	G	N3-C2-N2	-5.61	115.97	119.90
26	1H	340	A	O5'-P-OP2	5.61	117.43	110.70
26	1H	2469	A	N1-C6-N6	5.61	121.97	118.60
26	14	454	A	O5'-P-OP2	-5.61	100.65	105.70
26	14	2351	G	C6-C5-N7	-5.61	127.03	130.40
1	13	1027	C	OP1-P-O3'	5.61	117.53	105.20
26	1H	34	C	N1-C1'-C2'	5.61	121.29	114.00
26	1H	686	G	OP1-P-OP2	5.61	128.01	119.60
26	14	2627	G	C4-N9-C1'	5.61	133.79	126.50
26	1H	391	G	C2-N3-C4	-5.61	109.10	111.90
26	1H	678	C	OP1-P-O3'	5.61	117.53	105.20
26	1H	1258	C	OP2-P-O3'	5.60	117.53	105.20
26	1H	1400	G	N1-C6-O6	-5.60	116.54	119.90
26	1H	2060	A	N9-C4-C5	5.60	108.04	105.80
1	1G	275	G	N1-C6-O6	5.60	123.26	119.90
26	14	1320	C	N1-C2-O2	-5.60	115.54	118.90
26	1H	115	C	C5-C4-N4	-5.60	116.28	120.20
26	14	2334	G	N3-C4-N9	5.60	129.36	126.00
1	1G	817	C	C6-N1-C2	5.60	122.54	120.30
26	14	2638	G	P-O3'-C3'	5.60	126.42	119.70
26	1H	863	A	C8-N9-C4	5.60	108.04	105.80
26	1H	2577	A	N9-C4-C5	5.60	108.04	105.80
26	14	1407	C	C2-N3-C4	5.60	122.70	119.90
26	14	2607	G	N1-C6-O6	5.60	123.26	119.90
26	14	2818	G	N9-C4-C5	-5.60	103.16	105.40
1	13	858	G	C5-C6-O6	5.60	131.96	128.60
1	13	1413	A	C5-C6-N1	5.60	120.50	117.70
26	1H	622	G	C8-N9-C4	5.60	108.64	106.40
26	1H	1258	C	C2-N1-C1'	-5.60	112.64	118.80
23	2L	71	G	N3-C4-C5	5.60	131.40	128.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	2424	C	N3-C4-C5	5.60	124.14	121.90
26	1H	190	A	C5-C6-N6	-5.60	119.22	123.70
26	14	1663	C	N1-C2-O2	-5.60	115.54	118.90
1	13	656	C	C6-N1-C2	-5.59	118.06	120.30
1	13	803	G	C5-C6-O6	5.59	131.96	128.60
26	1H	967	C	O5'-P-OP1	5.59	117.41	110.70
26	1H	1229(A)	G	N1-C2-N3	5.59	127.26	123.90
26	1H	2700	C	C2-N3-C4	-5.59	117.10	119.90
26	14	2580	U	O5'-P-OP1	-5.59	100.67	105.70
26	14	2731	G	N1-C6-O6	5.59	123.26	119.90
29	29	80	GLU	N-CA-C	5.59	126.10	111.00
26	1H	263	C	C6-N1-C1'	-5.59	114.09	120.80
26	14	138	G	C5-C6-O6	-5.59	125.25	128.60
26	14	195	A	N1-C6-N6	5.59	121.95	118.60
26	14	465	G	OP1-P-OP2	-5.59	111.21	119.60
26	14	954	G	N1-C6-O6	-5.59	116.55	119.90
26	14	1971	A	N7-C8-N9	-5.59	111.00	113.80
22	1K	18	G	OP1-P-O3'	5.59	117.50	105.20
26	1H	910	A	O5'-P-OP1	5.59	117.41	110.70
26	14	365	C	C6-N1-C2	-5.59	118.06	120.30
26	14	762	U	C2-N1-C1'	5.59	124.41	117.70
26	14	2023	G	C8-N9-C4	-5.59	104.16	106.40
1	13	606	G	C2-N3-C4	5.59	114.69	111.90
26	1H	1410	G	N3-C4-C5	5.59	131.39	128.60
1	13	465	A	C5-N7-C8	5.59	106.69	103.90
26	1H	1156	A	N1-C6-N6	5.59	121.95	118.60
26	14	209	C	C6-N1-C1'	-5.59	114.09	120.80
26	14	631	A	OP1-P-O3'	5.59	117.49	105.20
26	14	2314	C	N3-C2-O2	-5.59	117.99	121.90
26	1H	906	G	C8-N9-C1'	5.58	134.26	127.00
26	1H	1521	G	N7-C8-N9	5.58	115.89	113.10
26	1H	1534	G	N3-C4-C5	-5.58	125.81	128.60
26	1H	1625	C	C5-C4-N4	5.58	124.11	120.20
26	1H	2426	A	O5'-P-OP2	-5.58	100.67	105.70
26	14	1334	G	C8-N9-C4	-5.58	104.17	106.40
26	14	2030	A	O5'-P-OP2	-5.58	100.67	105.70
1	13	1504	G	P-O3'-C3'	5.58	126.40	119.70
26	1H	247	G	N3-C2-N2	5.58	123.81	119.90
26	1H	952	G	C2-N3-C4	5.58	114.69	111.90
26	1H	1398	C	C6-N1-C2	5.58	122.53	120.30
26	1H	1702	G	C5-C6-N1	5.58	114.29	111.50
26	1H	2072	G	C5-C6-O6	-5.58	125.25	128.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
33	61	131	LYS	C-N-CD	-5.58	108.32	120.60
1	1G	995	C	C5-C6-N1	5.58	123.79	121.00
1	1G	1398	A	N1-C6-N6	-5.58	115.25	118.60
26	14	710	G	N1-C6-O6	5.58	123.25	119.90
26	14	1438	U	N3-C4-C5	-5.58	111.25	114.60
26	14	1922	G	N1-C6-O6	5.58	123.25	119.90
26	14	2012	G	N9-C4-C5	-5.58	103.17	105.40
26	14	2286	A	N1-C6-N6	5.58	121.95	118.60
1	13	449	C	C6-N1-C2	-5.58	118.07	120.30
26	1H	271(B)	G	C4-N9-C1'	5.58	133.76	126.50
26	1H	835	A	C5-C6-N1	5.58	120.49	117.70
26	14	733	G	N3-C4-C5	-5.58	125.81	128.60
26	1H	1851	U	N3-C2-O2	5.58	126.11	122.20
26	1H	1938	A	C6-C5-N7	-5.58	128.39	132.30
26	14	2555	U	N3-C2-O2	5.58	126.11	122.20
26	1H	658	C	N1-C2-O2	5.58	122.25	118.90
26	1H	792	G	N7-C8-N9	5.58	115.89	113.10
26	1H	1972	A	O5'-P-OP2	-5.58	100.68	105.70
26	1H	2443	C	N3-C4-N4	5.58	121.90	118.00
26	1H	2867	G	O4'-C1'-N9	5.58	112.66	108.20
26	14	583	G	C5-C6-N1	-5.58	108.71	111.50
26	14	1585	C	N3-C2-O2	-5.58	118.00	121.90
26	1H	245	G	C8-N9-C1'	-5.58	119.75	127.00
26	1H	850	C	C6-N1-C2	-5.58	118.07	120.30
26	1H	1698	A	N9-C1'-C2'	5.58	121.25	114.00
26	14	2607	G	C8-N9-C4	5.58	108.63	106.40
26	1H	102	G	OP1-P-O3'	5.58	117.47	105.20
26	1H	685	A	O4'-C1'-N9	5.58	112.66	108.20
26	1H	1573	G	OP1-P-O3'	-5.58	92.94	105.20
26	1H	1836	C	OP1-P-O3'	5.58	117.47	105.20
26	14	2066	C	N3-C4-C5	-5.58	119.67	121.90
26	14	2829	C	C5-C4-N4	-5.58	116.30	120.20
26	1H	1142	U	C5-C6-N1	5.57	125.49	122.70
26	1H	1842	G	OP2-P-O3'	5.57	117.46	105.20
26	1H	1917	U	OP1-P-O3'	5.57	117.46	105.20
26	1H	2466	C	N3-C4-C5	5.57	124.13	121.90
26	1H	2493	U	C6-N1-C2	-5.57	117.66	121.00
26	14	1307	A	N1-C2-N3	5.57	132.09	129.30
26	14	2087	G	N1-C6-O6	5.57	123.24	119.90
1	13	747	C	O5'-P-OP2	-5.57	100.69	105.70
26	1H	132	G	O5'-P-OP1	-5.57	100.69	105.70
26	1H	691	C	C2-N3-C4	-5.57	117.11	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	795	C	C2-N1-C1'	-5.57	112.67	118.80
26	1H	940	G	C6-N1-C2	-5.57	121.76	125.10
26	1H	1781	C	C6-N1-C1'	-5.57	114.12	120.80
26	1H	2848	G	N3-C2-N2	5.57	123.80	119.90
26	14	1938	A	C6-N1-C2	-5.57	115.26	118.60
26	14	2731	G	N7-C8-N9	5.57	115.88	113.10
26	14	1001	A	N1-C6-N6	-5.57	115.26	118.60
1	13	892	A	N1-C2-N3	5.57	132.08	129.30
26	1H	444	C	O5'-P-OP1	5.57	117.38	110.70
26	1H	627	A	N7-C8-N9	-5.57	111.02	113.80
26	1H	1948	G	C5-C6-O6	5.57	131.94	128.60
27	16	81	G	O4'-C1'-N9	5.57	112.65	108.20
26	1H	1229(A)	G	C6-C5-N7	-5.56	127.06	130.40
26	1H	708	C	C5-C6-N1	-5.56	118.22	121.00
26	1H	740	U	O5'-P-OP2	-5.56	100.69	105.70
1	1G	481	G	N3-C4-C5	-5.56	125.82	128.60
26	14	2331	G	C5-C6-O6	-5.56	125.26	128.60
1	13	1495	U	C6-N1-C2	-5.56	117.66	121.00
23	2K	17	C	C2-N3-C4	5.56	122.68	119.90
1	13	1199	U	N1-C2-N3	5.56	118.23	114.90
26	1H	411	G	C8-N9-C4	-5.56	104.18	106.40
26	1H	1393	A	O4'-C1'-N9	5.56	112.65	108.20
26	1H	2246	G	N3-C4-N9	5.56	129.34	126.00
26	14	150	C	C6-N1-C2	-5.56	118.08	120.30
26	14	1600	C	C5-C6-N1	-5.56	118.22	121.00
26	14	2253	G	C4-C5-N7	5.56	113.02	110.80
26	14	2377	A	C8-N9-C4	5.56	108.02	105.80
1	13	575	G	C5-C6-O6	5.56	131.94	128.60
23	2K	21	U	N3-C2-O2	-5.56	118.31	122.20
26	1H	120	U	C5-C4-O4	5.56	129.23	125.90
26	14	1282	U	N1-C2-N3	5.56	118.23	114.90
26	14	2037	G	N3-C4-C5	-5.56	125.82	128.60
26	1H	2331	G	N1-C6-O6	5.56	123.23	119.90
26	1H	1257	C	C4-C5-C6	5.55	120.18	117.40
26	1H	1204	A	C5-N7-C8	-5.55	101.12	103.90
26	14	1283	G	N3-C4-N9	5.55	129.33	126.00
26	14	1594	G	O5'-P-OP1	-5.55	100.70	105.70
27	16	98	G	C6-C5-N7	-5.55	127.07	130.40
26	14	475	U	C6-N1-C2	-5.55	117.67	121.00
26	1H	596	G	C5-C6-O6	-5.55	125.27	128.60
26	1H	928	G	O5'-P-OP2	5.55	117.36	110.70
26	1H	1156	A	N9-C4-C5	-5.55	103.58	105.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	2506	U	C5-C6-N1	5.55	125.47	122.70
1	1G	197	A	N7-C8-N9	5.55	116.58	113.80
26	14	580	C	N3-C4-N4	5.55	121.89	118.00
26	14	1356	G	O5'-P-OP1	-5.55	100.71	105.70
1	13	1414	U	OP2-P-O3'	5.55	117.40	105.20
26	1H	115	C	N3-C4-N4	5.55	121.88	118.00
26	1H	617	G	C5-N7-C8	5.55	107.07	104.30
26	1H	942	G	N9-C4-C5	5.55	107.62	105.40
26	1H	2247	A	C5-C6-N6	5.55	128.14	123.70
26	1H	2254	C	N1-C2-O2	-5.55	115.57	118.90
1	1G	1499	A	C8-N9-C4	5.55	108.02	105.80
26	14	945	A	N1-C2-N3	5.55	132.07	129.30
26	14	2056	G	N1-C2-N2	5.55	121.19	116.20
26	1H	925	C	O5'-P-OP1	5.54	117.35	110.70
26	1H	1625	C	N3-C4-N4	-5.54	114.12	118.00
26	14	1607	C	C6-N1-C1'	-5.54	114.15	120.80
26	1H	1284	A	O5'-P-OP2	-5.54	100.71	105.70
1	1G	230	G	C2-N3-C4	-5.54	109.13	111.90
26	1H	190	A	O5'-P-OP1	5.54	117.35	110.70
26	1H	1403	C	N3-C4-N4	-5.54	114.12	118.00
26	1H	2060	A	C8-N9-C1'	5.54	137.68	127.70
26	1H	2311	A	C5-N7-C8	-5.54	101.13	103.90
25	4L	21	C	C5-C6-N1	5.54	123.77	121.00
26	14	1314	C	C5-C6-N1	5.54	123.77	121.00
26	14	1769	G	N3-C4-C5	-5.54	125.83	128.60
1	13	584	G	N1-C2-N3	5.54	127.22	123.90
26	1H	239	U	C5-C6-N1	-5.54	119.93	122.70
26	1H	146	G	C8-N9-C4	5.54	108.62	106.40
26	1H	251	A	C4-C5-C6	5.54	119.77	117.00
26	1H	657	U	OP2-P-O3'	5.54	117.39	105.20
26	1H	1203	G	N3-C4-C5	-5.54	125.83	128.60
26	1H	1797	C	C4-C5-C6	5.54	120.17	117.40
26	1H	2308	G	C5-C6-N1	-5.54	108.73	111.50
26	1H	2490	G	C6-C5-N7	-5.54	127.08	130.40
26	1H	2554	U	C5-C6-N1	5.54	125.47	122.70
26	14	1029	A	C8-N9-C4	5.54	108.02	105.80
26	14	2233	U	N1-C2-N3	5.54	118.22	114.90
1	13	816	A	C5-C6-N6	5.54	128.13	123.70
1	13	1519	A	C4-C5-C6	5.54	119.77	117.00
26	1H	389	G	C8-N9-C1'	-5.54	119.80	127.00
26	1H	1344	G	C4-C5-N7	5.54	113.02	110.80
26	1H	2690	C	N1-C2-O2	-5.54	115.58	118.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	586	A	OP1-P-O3'	5.54	117.38	105.20
1	13	326	G	C4-C5-N7	-5.54	108.59	110.80
26	1H	1201	C	C5-C4-N4	-5.54	116.33	120.20
26	1H	2096	U	C5-C6-N1	5.54	125.47	122.70
26	1H	2257	U	N3-C2-O2	-5.54	118.33	122.20
26	1H	2416	C	OP2-P-O3'	5.54	117.38	105.20
27	16	56	G	N7-C8-N9	5.54	115.87	113.10
1	1G	345	C	P-O3'-C3'	5.54	126.34	119.70
26	14	569	U	OP1-P-O3'	5.54	117.38	105.20
1	13	1027	C	P-O3'-C3'	5.53	126.34	119.70
26	1H	803	U	C2-N3-C4	-5.53	123.68	127.00
26	1H	1158	C	C5-C6-N1	-5.53	118.23	121.00
26	1H	1695	G	N3-C4-N9	5.53	129.32	126.00
26	1H	1981	A	O4'-C1'-N9	-5.53	103.77	108.20
26	14	1931	U	OP1-P-OP2	-5.53	111.30	119.60
26	14	1899	G	N1-C6-O6	-5.53	116.58	119.90
26	14	2702	U	O4'-C1'-N1	5.53	112.62	108.20
53	K5	36	LEU	CA-CB-CG	5.53	128.02	115.30
26	1H	180	G	N3-C4-N9	5.53	129.32	126.00
26	1H	1705	G	C8-N9-C4	-5.53	104.19	106.40
26	1H	2544	G	C5-C6-O6	-5.53	125.28	128.60
26	14	1678	G	O5'-P-OP1	-5.53	100.72	105.70
26	1H	835	A	N9-C4-C5	5.53	108.01	105.80
26	1H	1990	C	N1-C2-N3	5.53	123.07	119.20
1	13	802	A	N1-C6-N6	5.53	121.92	118.60
26	1H	74	A	C6-C5-N7	-5.53	128.43	132.30
26	1H	1427	A	P-O3'-C3'	5.53	126.33	119.70
28	11	46	GLN	C-N-CA	-5.53	110.69	122.30
1	1G	17	U	N3-C4-O4	5.53	123.27	119.40
26	14	1443	G	C6-C5-N7	-5.53	127.08	130.40
26	14	1820	U	O5'-P-OP2	5.53	117.33	110.70
1	13	1354	C	C5-C6-N1	5.53	123.76	121.00
26	1H	429	A	N7-C8-N9	5.53	116.56	113.80
45	C5	39	VAL	N-CA-C	5.53	125.92	111.00
1	13	584	G	N3-C4-C5	-5.52	125.84	128.60
26	1H	795	C	N3-C4-N4	-5.52	114.14	118.00
26	1H	1191	G	OP1-P-OP2	5.52	127.88	119.60
26	14	667	U	N3-C4-O4	5.52	123.27	119.40
26	14	961	C	OP1-P-OP2	5.52	127.88	119.60
26	14	1253	A	C8-N9-C4	5.52	108.01	105.80
26	14	2575	C	N3-C4-N4	-5.52	114.14	118.00
26	1H	1502	C	OP2-P-O3'	5.52	117.34	105.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	2018	G	C6-C5-N7	-5.52	127.09	130.40
26	14	1309	G	N1-C6-O6	5.52	123.21	119.90
26	14	2351	G	C8-N9-C4	-5.52	104.19	106.40
26	14	2392	A	O5'-P-OP2	5.52	117.32	110.70
27	1J	71	C	C2-N1-C1'	5.52	124.87	118.80
26	1H	649	G	O5'-P-OP2	-5.52	100.73	105.70
26	14	602	G	C6-C5-N7	-5.52	127.09	130.40
26	14	1395	A	OP2-P-O3'	5.52	117.34	105.20
26	1H	705	A	C5-C6-N6	-5.52	119.29	123.70
26	1H	2041	U	O5'-P-OP1	-5.52	100.74	105.70
26	1H	58	G	OP2-P-O3'	5.51	117.33	105.20
26	1H	108	U	O5'-P-OP1	-5.51	100.74	105.70
26	1H	1010	A	OP2-P-O3'	5.51	117.33	105.20
26	1H	2357	U	OP2-P-O3'	5.51	117.33	105.20
26	14	2829	C	N3-C4-N4	5.51	121.86	118.00
26	1H	1254	A	C8-N9-C4	5.51	108.00	105.80
26	1H	2244	U	OP1-P-OP2	-5.51	111.33	119.60
26	14	265	A	C5-N7-C8	-5.51	101.14	103.90
26	14	1767	C	C2-N3-C4	-5.51	117.14	119.90
26	1H	912	C	O5'-P-OP2	5.51	117.31	110.70
26	1H	1574	C	N3-C4-C5	5.51	124.10	121.90
26	1H	2451	A	N9-C4-C5	5.51	108.00	105.80
26	14	1279	G	C5-C6-O6	5.51	131.91	128.60
26	14	2315	G	N3-C4-C5	-5.51	125.84	128.60
1	13	726	C	N1-C2-O2	5.51	122.21	118.90
26	1H	955	C	C6-N1-C2	-5.51	118.10	120.30
26	1H	1823	G	N3-C2-N2	-5.51	116.04	119.90
26	1H	2057	A	N7-C8-N9	-5.51	111.05	113.80
26	1H	2622	C	O5'-P-OP2	-5.51	100.74	105.70
26	1H	2745	C	C6-N1-C2	-5.51	118.10	120.30
26	14	756	C	N3-C4-C5	-5.51	119.70	121.90
26	14	1698	A	C5-C6-N1	-5.51	114.94	117.70
26	14	810	U	N1-C2-O2	-5.51	118.94	122.80
1	13	129	U	C6-N1-C1'	5.51	128.91	121.20
26	1H	1247	A	OP2-P-O3'	5.51	117.31	105.20
26	1H	2375	G	C8-N9-C4	5.51	108.60	106.40
26	1H	2487	G	N1-C6-O6	5.51	123.20	119.90
1	1G	615	C	C6-N1-C2	-5.51	118.10	120.30
26	1H	192	C	N3-C4-N4	5.50	121.85	118.00
26	1H	1604	C	N3-C4-C5	-5.50	119.70	121.90
26	14	940	G	N3-C4-C5	-5.50	125.85	128.60
26	14	1613	G	N7-C8-N9	-5.50	110.35	113.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	2596	U	N1-C2-N3	5.50	118.20	114.90
26	14	2829	C	N3-C2-O2	5.50	125.75	121.90
1	13	503	C	C6-N1-C2	-5.50	118.10	120.30
26	1H	767	U	O5'-P-OP2	-5.50	100.75	105.70
26	1H	791	C	OP1-P-O3'	-5.50	93.09	105.20
26	1H	1314	C	OP2-P-O3'	5.50	117.31	105.20
26	1H	2583	G	C8-N9-C4	-5.50	104.20	106.40
1	1G	1192	C	C5-C6-N1	5.50	123.75	121.00
26	14	791	C	N3-C4-C5	5.50	124.10	121.90
26	14	2249	U	C5-C6-N1	5.50	125.45	122.70
26	1H	94	G	N1-C6-O6	5.50	123.20	119.90
26	1H	2330	G	N1-C2-N3	5.50	127.20	123.90
1	1G	925	G	C8-N9-C4	5.50	108.60	106.40
26	1H	2075	U	C4-C5-C6	5.50	123.00	119.70
27	16	39	A	C8-N9-C4	-5.50	103.60	105.80
26	14	2592	G	N3-C4-C5	-5.50	125.85	128.60
1	13	1432	G	C6-C5-N7	-5.50	127.10	130.40
26	1H	610	C	C2-N3-C4	-5.50	117.15	119.90
26	1H	2039	C	N3-C4-C5	-5.50	119.70	121.90
26	14	1359	A	N1-C2-N3	-5.50	126.55	129.30
26	14	2565	A	N9-C4-C5	-5.50	103.60	105.80
1	13	233	C	C6-N1-C2	-5.50	118.10	120.30
26	1H	1805	U	O5'-P-OP1	-5.50	100.75	105.70
26	1H	2830	G	O5'-P-OP2	-5.50	100.75	105.70
26	14	1313	U	N3-C4-C5	-5.50	111.30	114.60
1	13	580	U	C2-N3-C4	-5.49	123.70	127.00
26	1H	674	G	OP1-P-OP2	-5.49	111.36	119.60
26	1H	814	C	O5'-P-OP2	-5.49	100.76	105.70
26	1H	970	C	O5'-P-OP1	-5.49	100.76	105.70
26	1H	1300	U	OP1-P-O3'	5.49	117.29	105.20
26	1H	1437	C	C2-N1-C1'	5.49	124.84	118.80
26	1H	2450	A	N1-C6-N6	-5.49	115.30	118.60
26	14	1399	C	OP2-P-O3'	5.49	117.29	105.20
26	14	2867	G	O4'-C1'-N9	5.49	112.59	108.20
1	13	525	C	C5-C6-N1	5.49	123.75	121.00
26	1H	511	U	OP1-P-OP2	-5.49	111.36	119.60
26	1H	1443	G	C8-N9-C1'	-5.49	119.86	127.00
1	1G	1281	U	C5-C6-N1	5.49	125.45	122.70
26	14	1566	A	N1-C2-N3	-5.49	126.55	129.30
26	14	1790	C	C5-C4-N4	-5.49	116.36	120.20
26	1H	404	C	N3-C2-O2	5.49	125.74	121.90
26	1H	738	G	O4'-C1'-N9	5.49	112.59	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	764	A	N1-C6-N6	5.49	121.89	118.60
26	1H	835	A	C6-N1-C2	-5.49	115.31	118.60
26	1H	972	G	C5-C6-O6	5.49	131.89	128.60
26	1H	1422	G	C8-N9-C4	-5.49	104.20	106.40
26	1H	1670	C	C2-N3-C4	-5.49	117.16	119.90
26	1H	2393	A	OP2-P-O3'	5.49	117.28	105.20
26	1H	2532	G	N1-C6-O6	5.49	123.19	119.90
26	14	803	U	N3-C2-O2	-5.49	118.36	122.20
26	14	1396	U	C4-C5-C6	5.49	122.99	119.70
26	14	2346	A	C5-N7-C8	-5.49	101.16	103.90
26	1H	1466	G	OP2-P-O3'	5.49	117.28	105.20
26	14	786	C	OP2-P-O3'	5.49	117.28	105.20
26	14	992	C	C6-N1-C2	-5.49	118.10	120.30
26	14	2258	C	OP1-P-O3'	5.49	117.28	105.20
1	13	318	G	N1-C6-O6	5.49	123.19	119.90
26	1H	389	G	N9-C4-C5	-5.49	103.20	105.40
26	14	961	C	O5'-P-OP2	-5.49	100.76	105.70
26	1H	528	A	C4-N9-C1'	-5.49	116.43	126.30
26	1H	845	G	OP1-P-O3'	5.49	117.27	105.20
26	1H	2711	A	P-O3'-C3'	5.49	126.28	119.70
1	1G	230	G	N3-C2-N2	-5.49	116.06	119.90
1	13	1530	G	C5-C6-O6	-5.48	125.31	128.60
26	1H	1816	G	O5'-P-OP2	5.48	117.28	110.70
26	1H	777	A	N1-C6-N6	-5.48	115.31	118.60
26	1H	2324	C	C6-N1-C1'	-5.48	114.22	120.80
26	1H	2434	A	C4-C5-C6	-5.48	114.26	117.00
26	1H	2435	A	C8-N9-C4	-5.48	103.61	105.80
26	14	866	A	C4-N9-C1'	5.48	136.16	126.30
26	1H	401	A	N1-C2-N3	5.48	132.04	129.30
1	1G	1390	U	C5-C4-O4	5.48	129.19	125.90
1	13	101	A	C8-N9-C4	-5.48	103.61	105.80
1	13	792	A	C3'-C2'-C1'	-5.48	97.12	101.50
1	13	1475	G	N7-C8-N9	5.48	115.84	113.10
26	1H	121	G	C4-C5-N7	5.48	112.99	110.80
26	1H	2713	A	N3-C4-N9	-5.48	123.02	127.40
26	14	840	C	N1-C2-O2	-5.48	115.61	118.90
26	14	2755	C	C5-C6-N1	5.48	123.74	121.00
1	13	537	G	O5'-P-OP1	-5.47	100.77	105.70
26	1H	2219	G	N1-C6-O6	5.47	123.18	119.90
26	1H	2497	A	O5'-P-OP2	-5.47	100.77	105.70
26	14	199	A	OP2-P-O3'	5.47	117.25	105.20
26	14	2318	G	N7-C8-N9	5.47	115.84	113.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
38	55	113	LEU	CA-CB-CG	5.47	127.89	115.30
26	1H	645	C	N1-C2-O2	5.47	122.18	118.90
26	1H	727	A	OP2-P-O3'	5.47	117.24	105.20
26	1H	1319	G	N3-C4-C5	-5.47	125.86	128.60
26	1H	1435	G	N1-C6-O6	-5.47	116.62	119.90
26	1H	2692	C	C6-N1-C2	-5.47	118.11	120.30
26	1H	2827	C	N3-C4-N4	5.47	121.83	118.00
27	16	15	A	O4'-C1'-N9	5.47	112.58	108.20
26	14	2066	C	O5'-P-OP2	5.47	117.27	110.70
26	1H	188	G	N1-C2-N2	-5.47	111.28	116.20
26	1H	688	U	C2-N3-C4	-5.47	123.72	127.00
26	1H	1299	G	N7-C8-N9	5.47	115.84	113.10
1	13	1348	U	C5-C4-O4	5.47	129.18	125.90
1	13	1385	G	N1-C6-O6	5.47	123.18	119.90
26	1H	429	A	N9-C4-C5	5.47	107.99	105.80
26	1H	736	C	N1-C2-O2	-5.47	115.62	118.90
26	1H	2586	C	N3-C2-O2	5.47	125.73	121.90
1	1G	115	G	C4-C5-N7	-5.47	108.61	110.80
26	14	747	U	N1-C2-O2	-5.47	118.97	122.80
26	1H	2564	A	OP1-P-O3'	5.47	117.23	105.20
1	1G	613	C	C6-N1-C2	-5.47	118.11	120.30
26	14	784	A	P-O3'-C3'	5.47	126.26	119.70
26	14	791	C	P-O3'-C3'	5.47	126.26	119.70
26	14	995	C	C2-N1-C1'	-5.47	112.78	118.80
26	1H	1579	A	C8-N9-C4	-5.47	103.61	105.80
26	1H	1942	C	C4-C5-C6	-5.47	114.67	117.40
1	1G	894	G	C4-C5-N7	5.47	112.99	110.80
26	14	762	U	O4'-C1'-N1	-5.47	103.83	108.20
26	14	783	A	C5-C6-N6	-5.47	119.33	123.70
26	1H	73	A	C2-N3-C4	5.46	113.33	110.60
26	1H	2251	G	OP1-P-O3'	5.46	117.22	105.20
1	1G	632	A	P-O3'-C3'	5.46	126.26	119.70
26	14	184	C	OP2-P-O3'	5.46	117.22	105.20
26	14	460	A	C5-C6-N6	-5.46	119.33	123.70
26	14	2041	U	OP2-P-O3'	5.46	117.22	105.20
26	14	2360	A	C2-N3-C4	-5.46	107.87	110.60
26	14	2618	G	C4-C5-N7	-5.46	108.61	110.80
26	1H	2072	G	C8-N9-C4	5.46	108.58	106.40
26	1H	2253	G	N3-C4-N9	-5.46	122.72	126.00
26	14	270(U)	C	C6-N1-C2	-5.46	118.11	120.30
26	14	396	G	N1-C6-O6	5.46	123.18	119.90
26	14	693	C	C2-N3-C4	-5.46	117.17	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	13	422	C	P-O3'-C3'	5.46	126.25	119.70
23	2K	77	A	O5'-P-OP1	-5.46	100.78	105.70
26	1H	410	G	O5'-P-OP1	-5.46	100.78	105.70
26	1H	963	U	O5'-P-OP2	5.46	117.25	110.70
26	1H	1305	C	OP1-P-OP2	-5.46	111.41	119.60
26	1H	1359	A	OP1-P-OP2	5.46	127.79	119.60
26	1H	1366	A	C2-N3-C4	-5.46	107.87	110.60
26	1H	2319	G	N3-C4-N9	5.46	129.28	126.00
26	1H	2586	C	OP1-P-O3'	5.46	117.22	105.20
27	16	115	G	N3-C4-N9	5.46	129.28	126.00
26	14	1600	C	O5'-P-OP2	-5.46	100.78	105.70
26	14	1642	G	O5'-P-OP1	-5.46	100.78	105.70
1	13	1498	U	N3-C2-O2	-5.46	118.38	122.20
26	1H	120	U	N1-C2-N3	5.46	118.18	114.90
26	1H	541	C	N3-C2-O2	-5.46	118.08	121.90
26	1H	1299	G	C5-C6-O6	-5.46	125.32	128.60
26	1H	2004	G	C2-N3-C4	-5.46	109.17	111.90
26	1H	2589	A	O5'-P-OP2	-5.46	100.79	105.70
26	1H	430	G	OP1-P-O3'	5.46	117.21	105.20
26	1H	812	C	N3-C2-O2	5.46	125.72	121.90
26	1H	1601	G	OP1-P-O3'	5.46	117.21	105.20
26	1H	2779	U	C5-C6-N1	-5.46	119.97	122.70
1	1G	975	A	O4'-C1'-N9	-5.46	103.83	108.20
26	14	1365	A	C8-N9-C4	-5.46	103.62	105.80
26	14	1928	A	C8-N9-C4	5.46	107.98	105.80
24	3K	76	A	O4'-C1'-N9	5.46	112.56	108.20
28	11	131	LEU	CA-CB-CG	5.46	127.85	115.30
1	1G	865	A	C8-N9-C4	-5.46	103.62	105.80
26	14	298	G	C5-N7-C8	-5.46	101.57	104.30
1	13	231	G	C8-N9-C4	-5.46	104.22	106.40
26	1H	600	G	C5-C6-N1	-5.46	108.77	111.50
26	1H	988	A	P-O3'-C3'	5.46	126.25	119.70
26	14	728	G	C8-N9-C4	5.46	108.58	106.40
26	1H	417	C	N3-C4-C5	5.45	124.08	121.90
26	1H	617	G	C5-C6-N1	5.45	114.23	111.50
26	1H	955	C	C4-C5-C6	5.45	120.13	117.40
26	1H	1192	G	N7-C8-N9	-5.45	110.37	113.10
26	1H	2360	A	C2-N3-C4	-5.45	107.87	110.60
26	1H	2606	C	OP1-P-O3'	5.45	117.20	105.20
26	14	2259	G	N1-C2-N2	5.45	121.11	116.20
1	13	1195	C	C6-N1-C2	-5.45	118.12	120.30
1	13	1519	A	C5-C6-N6	5.45	128.06	123.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	2022	U	N3-C2-O2	5.45	126.02	122.20
26	14	1355	G	N7-C8-N9	5.45	115.83	113.10
26	14	1831	G	N1-C2-N3	5.45	127.17	123.90
26	1H	1325	G	C5-C6-O6	-5.45	125.33	128.60
26	1H	2389	G	C8-N9-C4	-5.45	104.22	106.40
26	1H	2705	A	C4-C5-N7	5.45	113.42	110.70
26	14	93	C	C6-N1-C2	-5.45	118.12	120.30
1	13	557	G	N3-C4-N9	5.45	129.27	126.00
1	13	1065	U	P-O3'-C3'	5.45	126.24	119.70
26	1H	1241	A	C5-C6-N1	-5.45	114.98	117.70
26	14	127	A	OP1-P-O3'	5.45	117.19	105.20
26	14	2444	G	C4-C5-N7	-5.45	108.62	110.80
26	14	797	C	N3-C4-N4	5.45	121.81	118.00
26	14	802	A	C6-N1-C2	-5.45	115.33	118.60
26	14	1725	G	C4-N9-C1'	5.45	133.58	126.50
26	14	2443	C	C2-N3-C4	-5.45	117.18	119.90
1	1G	243	A	P-O3'-C3'	5.45	126.23	119.70
26	14	1463	C	N3-C4-C5	-5.45	119.72	121.90
26	14	1520	U	N3-C4-O4	-5.45	115.59	119.40
26	14	2618	G	C5-C6-O6	5.45	131.87	128.60
26	1H	1764	G	N1-C2-N3	5.44	127.17	123.90
1	13	14	U	C6-N1-C2	-5.44	117.73	121.00
26	1H	2508	G	C8-N9-C1'	5.44	134.07	127.00
1	1G	793	U	O4'-C1'-N1	5.44	112.55	108.20
26	14	632	A	O5'-P-OP2	5.44	117.23	110.70
1	13	581	G	C6-C5-N7	-5.44	127.14	130.40
1	13	945	G	OP1-P-O3'	5.44	117.17	105.20
1	1G	586	C	C6-N1-C2	-5.44	118.12	120.30
26	14	1327	C	C6-N1-C1'	5.44	127.33	120.80
26	1H	530	G	N3-C2-N2	5.44	123.71	119.90
26	1H	1616	A	OP1-P-O3'	5.44	117.17	105.20
26	1H	1688	U	OP2-P-O3'	5.44	117.17	105.20
26	1H	105	C	C6-N1-C2	-5.44	118.12	120.30
27	16	12	C	C4-C5-C6	5.44	120.12	117.40
26	14	459	U	O5'-P-OP2	-5.44	100.81	105.70
26	1H	389	G	N3-C4-N9	5.44	129.26	126.00
26	1H	1363	C	C2-N3-C4	-5.43	117.18	119.90
26	1H	2598	A	C8-N9-C4	5.43	107.97	105.80
1	13	503	C	C2-N1-C1'	5.43	124.78	118.80
1	13	963	G	N3-C4-C5	-5.43	125.88	128.60
26	1H	116	C	C6-N1-C1'	5.43	127.32	120.80
26	1H	713	G	N1-C6-O6	5.43	123.16	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	1577	C	C4-C5-C6	5.43	120.12	117.40
1	1G	950	U	C5-C6-N1	5.43	125.42	122.70
26	14	601	C	N1-C2-O2	-5.43	115.64	118.90
26	14	1266	G	C5-C6-N1	5.43	114.22	111.50
26	14	37	C	OP2-P-O3'	5.43	117.15	105.20
26	14	1024	G	C6-C5-N7	-5.43	127.14	130.40
26	1H	1189	A	C5-C6-N6	-5.43	119.36	123.70
1	1G	377	G	N9-C4-C5	-5.43	103.23	105.40
1	13	974	A	C6-C5-N7	-5.43	128.50	132.30
26	1H	1367	A	O5'-P-OP1	5.43	117.21	110.70
1	13	1335	C	C2-N1-C1'	-5.43	112.83	118.80
26	1H	2267	A	P-O3'-C3'	5.43	126.21	119.70
1	1G	132	C	N3-C4-C5	-5.43	119.73	121.90
1	1G	354	G	C8-N9-C1'	-5.43	119.95	127.00
1	1G	377	G	C6-C5-N7	-5.43	127.14	130.40
1	1G	858	G	C4-C5-N7	5.43	112.97	110.80
26	14	570	G	N1-C2-N2	-5.43	111.32	116.20
26	14	828	U	N3-C4-C5	-5.43	111.34	114.60
26	14	1836	C	OP1-P-O3'	5.43	117.14	105.20
1	13	335	C	C2-N1-C1'	-5.42	112.83	118.80
24	3K	34	G	C8-N9-C4	-5.42	104.23	106.40
26	1H	1777	U	N3-C4-C5	-5.42	111.34	114.60
26	1H	2439	A	OP1-P-OP2	5.42	127.74	119.60
26	1H	2871	C	C6-N1-C2	-5.42	118.13	120.30
1	1G	1112	C	C6-N1-C2	-5.42	118.13	120.30
26	14	2053	G	C8-N9-C4	5.42	108.57	106.40
26	14	2086	U	C5-C4-O4	5.42	129.15	125.90
26	1H	126	A	OP2-P-O3'	5.42	117.13	105.20
26	1H	818	G	C5-C6-O6	5.42	131.85	128.60
26	1H	955	C	C5-C4-N4	5.42	124.00	120.20
26	1H	1319	G	N7-C8-N9	5.42	115.81	113.10
26	1H	685	A	C6-C5-N7	-5.42	128.50	132.30
26	1H	1336	A	C8-N9-C4	-5.42	103.63	105.80
26	1H	2328	A	C4-C5-C6	5.42	119.71	117.00
26	14	330	A	C4-C5-N7	5.42	113.41	110.70
26	14	576	U	OP2-P-O3'	5.42	117.13	105.20
26	14	776	G	N1-C2-N2	5.42	121.08	116.20
1	13	114	U	O5'-P-OP2	-5.42	100.82	105.70
26	1H	1324	G	C5-C6-N1	-5.42	108.79	111.50
26	14	985	C	OP2-P-O3'	5.42	117.12	105.20
26	1H	85	G	O5'-P-OP1	5.42	117.20	110.70
26	1H	700	G	C5-C6-O6	-5.42	125.35	128.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	942	G	C4-C5-N7	-5.42	108.63	110.80
26	1H	1428	C	C2-N1-C1'	-5.42	112.84	118.80
1	13	19	C	C6-N1-C2	-5.42	118.13	120.30
26	1H	1786	A	N3-C4-C5	5.42	130.59	126.80
26	14	1639	U	N3-C4-O4	-5.42	115.61	119.40
26	14	2008	C	N1-C2-N3	5.42	122.99	119.20
26	14	2463	C	N3-C2-O2	5.42	125.69	121.90
26	1H	1385	G	N3-C4-C5	5.42	131.31	128.60
26	1H	2026	C	C4-C5-C6	5.42	120.11	117.40
26	14	2502	G	C5-N7-C8	-5.42	101.59	104.30
26	1H	1993	U	N1-C2-O2	-5.41	119.01	122.80
26	1H	2029	G	N7-C8-N9	5.41	115.81	113.10
1	1G	1511	G	C8-N9-C4	-5.41	104.23	106.40
26	14	1352	U	O5'-P-OP2	-5.41	100.83	105.70
26	14	1418	G	N1-C6-O6	5.41	123.15	119.90
26	14	1784	A	OP1-P-O3'	5.41	117.11	105.20
26	14	2232	U	O5'-P-OP2	-5.41	100.83	105.70
26	1H	206	U	N3-C4-O4	-5.41	115.61	119.40
26	1H	1269	A	N7-C8-N9	5.41	116.51	113.80
26	14	468	G	OP1-P-OP2	-5.41	111.48	119.60
1	13	101	A	C5-N7-C8	-5.41	101.19	103.90
1	13	805	C	C5-C6-N1	5.41	123.70	121.00
26	1H	1311	G	O5'-P-OP2	-5.41	100.83	105.70
26	1H	1643	G	N1-C6-O6	-5.41	116.65	119.90
26	1H	2228	G	C5-C6-O6	-5.41	125.35	128.60
26	14	389	G	C8-N9-C1'	-5.41	119.97	127.00
26	14	487	C	N1-C2-O2	-5.41	115.65	118.90
26	14	583	G	C8-N9-C4	-5.41	104.23	106.40
26	14	1336	A	C5-C6-N1	5.41	120.41	117.70
26	1H	2597	G	OP2-P-O3'	5.41	117.10	105.20
26	14	1782	C	N3-C2-O2	5.41	125.69	121.90
1	13	718	G	O5'-P-OP2	5.41	117.19	110.70
26	1H	1776	G	C5-C6-O6	-5.41	125.36	128.60
1	13	591	U	N3-C2-O2	-5.41	118.42	122.20
26	1H	715	G	N3-C4-C5	-5.41	125.90	128.60
26	1H	1520	U	OP2-P-O3'	5.41	117.09	105.20
26	14	570	G	C8-N9-C4	-5.41	104.24	106.40
26	14	1754	C	N1-C2-O2	5.41	122.14	118.90
26	14	2426	A	N1-C6-N6	5.41	121.84	118.60
26	1H	189	G	C2-N3-C4	-5.40	109.20	111.90
26	1H	446	G	C5-C6-N1	-5.40	108.80	111.50
26	1H	481	G	C5-C6-O6	-5.40	125.36	128.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	1G	1432	G	C6-C5-N7	-5.40	127.16	130.40
1	13	314	C	N3-C2-O2	-5.40	118.12	121.90
1	13	934	C	C2-N1-C1'	-5.40	112.86	118.80
26	14	728	G	N3-C2-N2	5.40	123.68	119.90
1	13	966	G	C5-C6-O6	-5.40	125.36	128.60
43	E8	23	LEU	CA-CB-CG	5.40	127.72	115.30
1	1G	817	C	C5-C6-N1	-5.40	118.30	121.00
26	14	1841	U	OP2-P-O3'	5.40	117.08	105.20
26	14	1979	C	O5'-P-OP2	-5.40	100.84	105.70
26	1H	843	G	OP2-P-O3'	5.40	117.08	105.20
27	1J	22	U	C6-N1-C2	-5.40	117.76	121.00
26	1H	196	A	O4'-C1'-N9	5.40	112.52	108.20
26	1H	671	C	C2-N3-C4	-5.40	117.20	119.90
26	1H	761	A	C5-N7-C8	-5.40	101.20	103.90
26	1H	2261	C	OP2-P-O3'	5.40	117.08	105.20
1	1G	121	C	N1-C2-O2	5.40	122.14	118.90
26	14	1616	A	N3-C4-C5	5.40	130.58	126.80
26	14	1695	G	C6-C5-N7	-5.40	127.16	130.40
26	14	1801	G	O5'-P-OP1	-5.40	100.84	105.70
26	14	1982	C	N1-C2-N3	-5.40	115.42	119.20
26	1H	2465	C	C4-C5-C6	5.40	120.10	117.40
26	1H	2509	G	N1-C6-O6	-5.40	116.66	119.90
26	14	1772	G	C5-C6-O6	-5.40	125.36	128.60
26	14	2447	G	O5'-P-OP2	-5.40	100.84	105.70
1	13	1469	G	C8-N9-C4	-5.39	104.24	106.40
26	1H	739	G	N1-C2-N3	-5.39	120.66	123.90
26	1H	2296	U	O5'-P-OP1	-5.39	100.84	105.70
26	14	2235	G	N3-C4-N9	5.39	129.24	126.00
26	1H	673	C	N1-C2-O2	-5.39	115.67	118.90
26	1H	951	C	N3-C4-N4	-5.39	114.22	118.00
26	1H	1475	G	N3-C2-N2	-5.39	116.12	119.90
26	1H	1695	G	N3-C4-C5	-5.39	125.90	128.60
26	1H	1924	C	N1-C2-O2	5.39	122.14	118.90
26	1H	2367	G	N3-C2-N2	-5.39	116.13	119.90
1	1G	17	U	C5-C4-O4	-5.39	122.67	125.90
26	14	1826	G	C4-C5-N7	-5.39	108.64	110.80
26	1H	799	G	C5-N7-C8	5.39	107.00	104.30
26	1H	945	A	O5'-P-OP2	-5.39	100.85	105.70
26	14	1770	G	N7-C8-N9	5.39	115.80	113.10
1	13	950	U	OP1-P-O3'	5.39	117.06	105.20
26	1H	749	C	C6-N1-C2	5.39	122.46	120.30
1	1G	1414	U	C5-C4-O4	5.39	129.13	125.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	133	C	C6-N1-C2	5.39	122.45	120.30
26	14	508	G	OP1-P-OP2	5.39	127.68	119.60
1	13	1522	U	C5-C6-N1	-5.39	120.01	122.70
26	14	540	G	N3-C4-N9	-5.39	122.77	126.00
26	14	1385	G	N3-C4-C5	5.39	131.29	128.60
26	1H	700	G	N1-C6-O6	5.39	123.13	119.90
26	1H	1231	G	N1-C6-O6	5.39	123.13	119.90
26	1H	2378	A	N9-C4-C5	-5.39	103.65	105.80
1	1G	1314	C	C6-N1-C2	-5.39	118.14	120.30
26	14	2440	C	OP1-P-O3'	5.39	117.05	105.20
23	2K	40	C	C5-C6-N1	5.38	123.69	121.00
26	1H	825	C	N3-C2-O2	5.38	125.67	121.90
26	14	1300	U	O5'-P-OP1	5.38	117.16	110.70
26	14	2581	G	C4-N9-C1'	5.38	133.50	126.50
26	1H	674	G	O5'-P-OP2	5.38	117.16	110.70
26	1H	1594	G	OP1-P-O3'	5.38	117.04	105.20
26	14	461	C	OP1-P-OP2	5.38	127.67	119.60
26	1H	607	U	C5-C4-O4	5.38	129.13	125.90
26	1H	734	A	OP1-P-OP2	5.38	127.67	119.60
26	1H	1670	C	C6-N1-C2	5.38	122.45	120.30
26	1H	2457	U	N1-C2-O2	-5.38	119.03	122.80
27	16	5	C	C5-C4-N4	-5.38	116.43	120.20
27	16	29	A	OP1-P-OP2	-5.38	111.53	119.60
1	1G	1070	U	N3-C2-O2	-5.38	118.43	122.20
1	1G	1228	C	O5'-P-OP2	-5.38	100.86	105.70
1	1G	1469	G	N1-C6-O6	5.38	123.13	119.90
26	14	102	G	O5'-P-OP1	-5.38	100.86	105.70
26	14	2066	C	C6-N1-C2	-5.38	118.15	120.30
26	1H	2553	G	OP1-P-O3'	5.38	117.03	105.20
25	4K	18	G	N3-C4-N9	-5.38	122.77	126.00
26	1H	2378	A	O5'-P-OP2	5.38	117.16	110.70
1	1G	586	C	C5-C6-N1	5.38	123.69	121.00
26	14	1814	G	C8-N9-C4	-5.38	104.25	106.40
26	1H	121	G	C5-C6-N1	5.38	114.19	111.50
26	1H	447	A	N1-C6-N6	-5.38	115.37	118.60
26	1H	937	U	C6-N1-C2	5.38	124.23	121.00
26	1H	1123	C	C2-N3-C4	-5.38	117.21	119.90
26	1H	1811	G	N9-C4-C5	5.38	107.55	105.40
26	1H	2232	U	O5'-P-OP2	-5.38	100.86	105.70
26	1H	2392	A	C6-N1-C2	5.38	121.83	118.60
46	H8	117	LEU	CA-CB-CG	5.38	127.67	115.30
26	14	494	G	C5-C6-N1	-5.38	108.81	111.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	1929	G	OP1-P-OP2	5.38	127.67	119.60
26	1H	957	A	C6-C5-N7	-5.38	128.54	132.30
27	16	6	C	C6-N1-C2	5.38	122.45	120.30
26	14	116	C	C6-N1-C2	-5.38	118.15	120.30
27	1J	60	C	C6-N1-C2	-5.38	118.15	120.30
1	13	693	G	N1-C6-O6	-5.37	116.68	119.90
1	13	1208	C	N1-C2-O2	-5.37	115.68	118.90
26	1H	1210	A	N3-C4-C5	5.37	130.56	126.80
26	1H	1406	U	OP1-P-O3'	5.37	117.02	105.20
26	1H	2096	U	N3-C2-O2	-5.37	118.44	122.20
26	14	932	G	N3-C4-N9	-5.37	122.78	126.00
26	1H	599	G	N1-C6-O6	-5.37	116.68	119.90
26	1H	2330	G	C5-N7-C8	-5.37	101.61	104.30
26	1H	2615	U	N1-C2-O2	5.37	126.56	122.80
55	Q8	52	LYS	C-N-CA	5.37	144.56	122.00
26	14	1308	A	N1-C2-N3	5.37	131.99	129.30
26	14	1849	G	N9-C4-C5	5.37	107.55	105.40
26	14	2595	G	N3-C4-C5	5.37	131.28	128.60
26	1H	52	A	O5'-P-OP2	-5.37	100.87	105.70
26	1H	866	A	C6-N1-C2	5.37	121.82	118.60
1	1G	400	C	N1-C2-O2	-5.37	115.68	118.90
1	1G	1285	A	P-O3'-C3'	5.37	126.14	119.70
24	1L	74	C	N3-C4-N4	-5.37	114.24	118.00
26	14	209	C	OP2-P-O3'	5.37	117.01	105.20
26	14	1372	U	C4-C5-C6	5.37	122.92	119.70
26	14	1443	G	C4-C5-C6	5.37	122.02	118.80
26	14	2573	C	C2-N1-C1'	5.37	124.70	118.80
1	13	1304	G	N3-C4-C5	-5.37	125.92	128.60
26	1H	859	G	N3-C4-N9	-5.37	122.78	126.00
26	1H	1394	U	C5-C6-N1	5.37	125.38	122.70
55	Q8	60	LEU	CB-CG-CD2	5.37	120.12	111.00
26	14	1657	C	OP1-P-O3'	5.37	117.01	105.20
26	14	1698	A	N7-C8-N9	5.37	116.48	113.80
1	13	465	A	N7-C8-N9	-5.37	111.12	113.80
26	1H	590	A	C8-N9-C4	-5.37	103.65	105.80
26	1H	2082	A	N7-C8-N9	-5.37	111.12	113.80
26	14	248	G	C6-C5-N7	-5.37	127.18	130.40
26	14	2381	C	C6-N1-C2	5.37	122.45	120.30
1	13	934	C	C6-N1-C1'	5.36	127.24	120.80
26	1H	638	G	OP1-P-OP2	5.36	127.65	119.60
26	1H	690	G	C5-C6-O6	-5.36	125.38	128.60
26	1H	2234	G	O5'-P-OP1	5.36	117.14	110.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	2441	C	OP1-P-O3'	5.36	117.00	105.20
26	14	2611	U	C5-C6-N1	5.36	125.38	122.70
26	1H	2783	G	C6-C5-N7	-5.36	127.18	130.40
26	1H	853	G	C8-N9-C4	5.36	108.55	106.40
26	14	1898	U	C6-N1-C2	-5.36	117.78	121.00
26	1H	971	C	C4-C5-C6	5.36	120.08	117.40
26	1H	137(A)	G	O4'-C1'-N9	5.36	112.49	108.20
26	1H	617	G	N1-C6-O6	-5.36	116.69	119.90
26	1H	735	A	C5-C6-N6	-5.36	119.41	123.70
27	16	101	A	C8-N9-C4	5.36	107.94	105.80
26	14	1930	G	N9-C4-C5	5.36	107.54	105.40
1	13	1299	A	N1-C6-N6	5.36	121.81	118.60
1	13	1433	A	O5'-P-OP1	-5.36	100.88	105.70
26	1H	982	C	OP1-P-O3'	5.36	116.98	105.20
26	1H	1024	G	N3-C4-C5	-5.36	125.92	128.60
26	1H	1608	A	O5'-P-OP1	-5.36	100.88	105.70
1	1G	266	G	O4'-C1'-N9	-5.36	103.92	108.20
26	14	182	A	OP2-P-O3'	5.36	116.98	105.20
26	14	1304	C	N3-C4-C5	5.36	124.04	121.90
26	1H	2281	C	C5-C4-N4	-5.35	116.45	120.20
26	14	1270	C	OP2-P-O3'	5.35	116.98	105.20
1	13	974	A	O4'-C1'-N9	5.35	112.48	108.20
26	1H	701	G	N1-C6-O6	5.35	123.11	119.90
26	1H	701	G	C5-C6-N1	-5.35	108.82	111.50
26	1H	757	U	OP1-P-OP2	5.35	127.63	119.60
26	1H	2014	A	N1-C6-N6	5.35	121.81	118.60
26	1H	2036	C	N3-C4-C5	-5.35	119.76	121.90
26	14	1286	A	OP2-P-O3'	5.35	116.98	105.20
26	14	1321	A	N1-C6-N6	5.35	121.81	118.60
26	14	1696	G	O5'-P-OP2	-5.35	100.88	105.70
1	13	809	G	O5'-P-OP1	-5.35	100.88	105.70
26	1H	1284	A	C5-C6-N1	-5.35	115.02	117.70
26	14	2329	G	N3-C4-C5	-5.35	125.92	128.60
26	1H	113	G	C6-C5-N7	-5.35	127.19	130.40
26	1H	463	G	N9-C4-C5	-5.35	103.26	105.40
26	1H	1610	A	C5-C6-N1	-5.35	115.03	117.70
1	1G	301	G	N1-C6-O6	5.35	123.11	119.90
1	1G	667	G	C5-C6-O6	-5.35	125.39	128.60
26	14	288	C	C2-N1-C1'	5.35	124.68	118.80
26	14	659	C	C6-N1-C2	5.35	122.44	120.30
26	14	852	G	N1-C6-O6	-5.35	116.69	119.90
33	69	131	LYS	C-N-CD	-5.35	108.83	120.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	13	219	C	C5-C6-N1	5.35	123.67	121.00
1	13	1081	G	OP1-P-O3'	5.35	116.96	105.20
26	1H	1300	U	C5-C4-O4	5.35	129.11	125.90
1	1G	506	G	O5'-P-OP2	5.35	117.12	110.70
26	14	1419	A	C8-N9-C4	5.35	107.94	105.80
26	1H	617	G	O5'-P-OP2	5.34	117.11	110.70
26	1H	1366	A	N1-C2-N3	5.34	131.97	129.30
26	1H	1391	U	C2-N1-C1'	5.34	124.11	117.70
26	1H	1780	A	N1-C2-N3	5.34	131.97	129.30
26	1H	2587	A	OP2-P-O3'	5.34	116.96	105.20
26	1H	2698	U	OP1-P-OP2	5.34	127.62	119.60
1	1G	293	G	N1-C6-O6	5.34	123.11	119.90
26	14	2080	G	N1-C6-O6	-5.34	116.69	119.90
26	14	2253	G	O5'-P-OP1	5.34	117.11	110.70
26	14	2283	C	N3-C4-N4	5.34	121.74	118.00
26	14	2731	G	C4-C5-C6	5.34	122.01	118.80
23	2K	22	A	C8-N9-C4	-5.34	103.66	105.80
26	1H	1437	C	N3-C2-O2	-5.34	118.16	121.90
34	15	2	LYS	N-CA-C	-5.34	96.57	111.00
26	1H	729	G	OP2-P-O3'	5.34	116.95	105.20
26	1H	2513	G	N3-C4-N9	5.34	129.21	126.00
26	14	733	G	O5'-P-OP2	-5.34	100.89	105.70
26	14	1586	A	C8-N9-C4	-5.34	103.66	105.80
26	14	2080	G	O5'-P-OP2	-5.34	100.89	105.70
26	14	2406	U	OP1-P-O3'	5.34	116.95	105.20
26	1H	1626	G	N1-C6-O6	5.34	123.10	119.90
26	1H	1690	A	N1-C6-N6	-5.34	115.40	118.60
1	1G	596	C	C6-N1-C2	5.34	122.44	120.30
26	1H	1303	G	N1-C6-O6	-5.34	116.70	119.90
26	1H	1779	U	C6-N1-C2	5.34	124.20	121.00
26	14	1664	A	O4'-C1'-N9	-5.34	103.93	108.20
26	14	2278	A	N9-C4-C5	5.34	107.94	105.80
26	1H	216	A	C8-N9-C4	5.34	107.94	105.80
26	1H	790	C	N3-C2-O2	5.34	125.64	121.90
26	1H	1147	C	C6-N1-C2	5.34	122.44	120.30
26	1H	81	G	C5-C6-O6	5.33	131.80	128.60
26	1H	1942	C	C6-N1-C2	-5.33	118.17	120.30
26	14	566	U	C5-C6-N1	-5.33	120.03	122.70
26	14	1770	G	O5'-P-OP2	5.33	117.10	110.70
26	1H	24	G	C8-N9-C4	-5.33	104.27	106.40
26	1H	946	G	C4-C5-N7	-5.33	108.67	110.80
26	1H	1123	C	C4-C5-C6	5.33	120.07	117.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	1900	A	C5-C6-N6	-5.33	119.43	123.70
26	1H	2068	U	OP1-P-O3'	5.33	116.94	105.20
26	14	2510	C	OP1-P-OP2	5.33	127.60	119.60
1	13	477	G	N3-C4-C5	5.33	131.27	128.60
26	1H	245	G	C4-N9-C1'	5.33	133.43	126.50
26	1H	906	G	N3-C4-N9	-5.33	122.80	126.00
26	1H	1785	A	C5-C6-N1	-5.33	115.03	117.70
26	1H	1977	A	N9-C4-C5	-5.33	103.67	105.80
26	1H	2604	U	N3-C4-C5	-5.33	111.40	114.60
26	1H	2616	C	C6-N1-C2	-5.33	118.17	120.30
26	1H	80	G	C8-N9-C4	-5.33	104.27	106.40
26	1H	574	C	N1-C2-N3	-5.33	115.47	119.20
26	14	761	A	C4-C5-C6	-5.33	114.34	117.00
26	14	1386	C	C2-N3-C4	5.33	122.56	119.90
26	14	2004	G	N3-C2-N2	-5.33	116.17	119.90
26	1H	455	C	C4-C5-C6	-5.33	114.74	117.40
26	1H	514	A	OP1-P-O3'	5.33	116.92	105.20
26	1H	526	A	C8-N9-C4	-5.33	103.67	105.80
26	1H	1209	G	N3-C2-N2	-5.33	116.17	119.90
26	1H	1974	C	O5'-P-OP2	-5.33	100.91	105.70
26	1H	2561	A	OP1-P-OP2	5.33	127.59	119.60
26	14	679	C	N1-C2-O2	-5.33	115.70	118.90
26	14	1474	C	C2-N3-C4	5.33	122.56	119.90
26	14	1528	A	C8-N9-C4	-5.33	103.67	105.80
26	14	1926	U	N3-C2-O2	-5.33	118.47	122.20
26	1H	728	G	OP2-P-O3'	5.32	116.91	105.20
26	1H	763	G	N3-C4-N9	5.32	129.19	126.00
26	1H	2571	C	C6-N1-C2	-5.32	118.17	120.30
26	14	756	C	C6-N1-C2	-5.32	118.17	120.30
26	14	2712(A)	A	C8-N9-C4	5.32	107.93	105.80
1	13	880	C	C6-N1-C2	5.32	122.43	120.30
26	1H	391	G	N9-C4-C5	-5.32	103.27	105.40
26	1H	799	G	OP1-P-OP2	-5.32	111.62	119.60
26	1H	2267	A	OP2-P-O3'	-5.32	93.49	105.20
26	1H	2373	G	C6-N1-C2	-5.32	121.91	125.10
26	14	2615	U	C2-N1-C1'	5.32	124.09	117.70
1	13	532	A	O4'-C1'-N9	5.32	112.46	108.20
1	13	1513	A	OP2-P-O3'	5.32	116.90	105.20
26	1H	471	A	C5-C6-N1	-5.32	115.04	117.70
26	1H	631	A	C8-N9-C4	5.32	107.93	105.80
26	1H	939	G	C5-C6-O6	5.32	131.79	128.60
26	1H	1257	C	N1-C2-O2	-5.32	115.71	118.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	1407	C	OP1-P-O3'	5.32	116.91	105.20
27	16	47	C	N3-C2-O2	5.32	125.62	121.90
1	13	1205	U	C5-C4-O4	5.32	129.09	125.90
26	1H	2295	C	C5-C6-N1	5.32	123.66	121.00
1	13	1505	G	O5'-P-OP2	5.32	117.08	110.70
26	1H	271(B)	G	C8-N9-C1'	-5.32	120.09	127.00
1	1G	108	G	C6-C5-N7	-5.32	127.21	130.40
26	14	1355	G	C8-N9-C4	-5.32	104.27	106.40
26	14	1657	C	N3-C2-O2	-5.32	118.18	121.90
26	14	1956	U	O5'-P-OP2	-5.32	100.91	105.70
26	14	2817	G	OP1-P-OP2	5.32	127.58	119.60
26	1H	1310	G	O5'-P-OP1	-5.32	100.92	105.70
1	1G	293	G	C6-C5-N7	-5.32	127.21	130.40
1	1G	913	A	OP2-P-O3'	5.32	116.89	105.20
26	14	2767	C	C6-N1-C2	-5.32	118.17	120.30
1	13	1299	A	C8-N9-C4	-5.31	103.67	105.80
26	1H	27	G	O4'-C1'-N9	5.31	112.45	108.20
26	14	1339	G	O5'-P-OP1	-5.31	100.92	105.70
26	1H	1373	A	O5'-P-OP2	-5.31	100.92	105.70
26	14	810	U	OP1-P-O3'	5.31	116.89	105.20
26	14	1558	A	P-O3'-C3'	5.31	126.07	119.70
26	14	1784	A	N1-C2-N3	5.31	131.96	129.30
27	1J	75	G	O5'-P-OP2	-5.31	100.92	105.70
1	13	129	U	C5-C4-O4	5.31	129.09	125.90
26	1H	1647	G	O5'-P-OP2	5.31	117.07	110.70
26	1H	2424	C	O5'-P-OP1	-5.31	100.92	105.70
1	13	1329	A	C4-C5-N7	5.31	113.36	110.70
26	1H	1298	C	C5-C6-N1	5.31	123.66	121.00
26	1H	1356	G	N3-C4-C5	5.31	131.25	128.60
26	1H	1962	C	N3-C2-O2	5.31	125.62	121.90
26	14	788	A	N1-C6-N6	5.31	121.79	118.60
26	14	2367	G	N7-C8-N9	5.31	115.75	113.10
26	1H	217	G	C4-C5-N7	-5.31	108.68	110.80
26	1H	528	A	N7-C8-N9	5.31	116.45	113.80
26	1H	1785	A	OP2-P-O3'	5.31	116.88	105.20
26	1H	1814	G	C6-N1-C2	-5.31	121.92	125.10
26	1H	2822	G	C4-C5-N7	5.31	112.92	110.80
26	14	951	C	OP1-P-O3'	5.31	116.88	105.20
26	14	963	U	O5'-P-OP2	5.31	117.07	110.70
26	14	1402	C	N3-C4-C5	-5.31	119.78	121.90
26	14	2844	G	C5-C6-O6	-5.31	125.42	128.60
27	1J	71	C	C6-N1-C2	-5.31	118.18	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	677	A	N1-C6-N6	-5.31	115.42	118.60
26	1H	1334	G	C8-N9-C4	-5.31	104.28	106.40
26	14	2042	A	C8-N9-C4	5.31	107.92	105.80
27	1J	81	G	C4-C5-N7	5.31	112.92	110.80
26	1H	2477	C	C6-N1-C2	-5.30	118.18	120.30
26	14	1367	A	C5-C6-N6	-5.30	119.46	123.70
23	2K	48	U	OP2-P-O3'	5.30	116.87	105.20
26	1H	1973	G	N3-C4-C5	-5.30	125.95	128.60
26	1H	2439	A	C2-N3-C4	-5.30	107.95	110.60
26	1H	309	G	C8-N9-C4	-5.30	104.28	106.40
26	1H	552	G	O5'-P-OP1	5.30	117.06	110.70
26	1H	815	C	N3-C4-C5	5.30	124.02	121.90
26	1H	945	A	N9-C1'-C2'	5.30	120.89	114.00
28	19	43	ARG	NE-CZ-NH2	-5.30	117.65	120.30
26	1H	389	G	C8-N9-C4	5.30	108.52	106.40
26	1H	590	A	N1-C2-N3	5.30	131.95	129.30
26	1H	680	G	O5'-P-OP1	-5.30	100.93	105.70
26	1H	1546	C	C5-C6-N1	5.30	123.65	121.00
26	1H	2766	G	N1-C6-O6	5.30	123.08	119.90
42	95	18	LEU	CA-CB-CG	5.30	127.49	115.30
1	13	901	A	C5-N7-C8	-5.30	101.25	103.90
26	1H	2354	G	OP1-P-O3'	5.30	116.86	105.20
27	16	41	U	O5'-P-OP1	-5.30	100.93	105.70
26	1H	480	A	N1-C6-N6	5.30	121.78	118.60
26	1H	2434	A	OP2-P-O3'	5.30	116.85	105.20
1	1G	943	U	C2-N1-C1'	5.30	124.06	117.70
26	14	2387	U	OP2-P-O3'	5.30	116.85	105.20
26	14	2440	C	C2-N1-C1'	-5.30	112.97	118.80
26	1H	2424	C	C4-C5-C6	-5.29	114.75	117.40
1	1G	1062	U	O5'-P-OP2	-5.29	100.94	105.70
26	14	828	U	OP2-P-O3'	5.29	116.85	105.20
26	1H	1198	U	C5-C6-N1	-5.29	120.05	122.70
26	1H	1787	A	C2-N3-C4	-5.29	107.95	110.60
26	1H	2381	C	C5-C4-N4	-5.29	116.50	120.20
26	1H	2867	G	O5'-P-OP1	-5.29	100.94	105.70
26	14	2509	G	N1-C6-O6	5.29	123.08	119.90
1	13	108	G	C4-C5-N7	5.29	112.92	110.80
26	1H	1621	U	C5-C4-O4	-5.29	122.72	125.90
26	1H	1705	G	OP1-P-OP2	-5.29	111.66	119.60
26	14	1821	A	C6-C5-N7	-5.29	128.60	132.30
26	14	2873	A	N9-C1'-C2'	5.29	120.88	114.00
26	1H	685	A	C4-C5-N7	5.29	113.34	110.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	764	A	OP1-P-O3'	5.29	116.84	105.20
26	1H	2299	G	N7-C8-N9	5.29	115.75	113.10
26	1H	2435	A	N7-C8-N9	5.29	116.44	113.80
27	16	98	G	N9-C4-C5	-5.29	103.28	105.40
26	14	140	A	C5-C6-N6	-5.29	119.47	123.70
26	14	1908	C	N3-C4-C5	-5.29	119.78	121.90
26	14	2084	C	N1-C2-O2	-5.29	115.73	118.90
26	14	2610	C	N1-C2-N3	-5.29	115.50	119.20
23	2K	62	C	C6-N1-C2	-5.29	118.19	120.30
26	1H	645	C	C5-C6-N1	5.29	123.64	121.00
26	1H	1959	G	C2-N3-C4	5.29	114.54	111.90
26	1H	2253	G	N9-C4-C5	5.29	107.52	105.40
26	14	127	A	O5'-P-OP2	-5.29	100.94	105.70
26	14	503	A	N9-C4-C5	5.29	107.92	105.80
26	14	866	A	N9-C4-C5	-5.29	103.69	105.80
26	1H	1382	G	C6-C5-N7	-5.29	127.23	130.40
26	1H	2488	A	N1-C2-N3	5.29	131.94	129.30
1	13	644	G	O5'-P-OP2	-5.29	100.94	105.70
26	1H	58	G	N9-C4-C5	5.29	107.51	105.40
26	1H	321	G	N9-C4-C5	-5.29	103.29	105.40
26	1H	844	C	N3-C4-C5	-5.29	119.79	121.90
26	1H	2511	U	N3-C2-O2	-5.29	118.50	122.20
26	1H	2711	A	OP2-P-O3'	-5.29	93.57	105.20
26	14	2585	U	OP1-P-O3'	5.29	116.83	105.20
1	13	570	G	C2-N3-C4	5.28	114.54	111.90
26	1H	448	U	C4-C5-C6	5.28	122.87	119.70
26	1H	683	C	C5-C4-N4	-5.28	116.50	120.20
26	1H	828	U	C6-N1-C1'	-5.28	113.80	121.20
49	K8	59	ARG	NE-CZ-NH2	5.28	122.94	120.30
26	14	1142	U	C6-N1-C1'	-5.28	113.80	121.20
26	14	1290	C	OP1-P-OP2	5.28	127.52	119.60
26	14	1292	U	C5-C4-O4	-5.28	122.73	125.90
26	14	1598	C	C2-N1-C1'	5.28	124.61	118.80
26	14	1786	A	N3-C4-C5	5.28	130.50	126.80
26	1H	1294	U	O5'-P-OP1	-5.28	100.95	105.70
26	14	1673	U	O5'-P-OP2	5.28	117.04	110.70
26	14	1904	G	O5'-P-OP2	-5.28	100.95	105.70
1	13	542	G	C5-N7-C8	-5.28	101.66	104.30
26	1H	335	C	C2-N3-C4	5.28	122.54	119.90
26	1H	1901	A	OP1-P-O3'	5.28	116.82	105.20
26	1H	2199	A	C8-N9-C4	-5.28	103.69	105.80
26	1H	2665	A	O4'-C1'-N9	5.28	112.42	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	1G	894	G	N1-C6-O6	5.28	123.07	119.90
26	14	1271	G	C5-C6-O6	5.28	131.77	128.60
26	1H	2532	G	C6-C5-N7	-5.28	127.23	130.40
1	1G	932	C	N3-C2-O2	-5.28	118.20	121.90
26	14	1313	U	OP1-P-O3'	5.28	116.81	105.20
1	13	53	A	N1-C6-N6	5.28	121.77	118.60
26	1H	678	C	P-O3'-C3'	-5.28	113.37	119.70
26	1H	713	G	C5-C6-N1	-5.28	108.86	111.50
26	1H	1325	G	N1-C6-O6	5.28	123.07	119.90
27	16	19	G	N3-C4-C5	5.28	131.24	128.60
26	14	945	A	C5-C6-N1	-5.28	115.06	117.70
26	14	1349	A	N1-C6-N6	5.28	121.77	118.60
1	13	275	G	N3-C4-N9	5.28	129.17	126.00
26	1H	629	G	C2-N3-C4	-5.28	109.26	111.90
26	1H	1905	C	OP1-P-OP2	-5.28	111.69	119.60
26	1H	2506	U	C2-N1-C1'	5.28	124.03	117.70
1	13	1491	G	OP2-P-O3'	5.27	116.80	105.20
26	1H	663	G	C4-N9-C1'	5.27	133.36	126.50
26	1H	682	G	C8-N9-C4	5.27	108.51	106.40
26	14	130	C	C5-C4-N4	-5.27	116.51	120.20
26	1H	354	G	C5-C6-O6	-5.27	125.44	128.60
26	1H	410	G	O5'-P-OP2	5.27	117.03	110.70
26	1H	1908	C	O5'-P-OP2	-5.27	100.96	105.70
26	1H	2363	C	OP2-P-O3'	5.27	116.80	105.20
26	14	768	G	O5'-P-OP2	-5.27	100.95	105.70
26	14	2230	G	C5-C6-O6	5.27	131.76	128.60
26	1H	533	G	C4-C5-N7	-5.27	108.69	110.80
26	1H	1790	C	C2-N3-C4	-5.27	117.26	119.90
26	14	661	C	N3-C4-C5	-5.27	119.79	121.90
1	13	701	C	N1-C2-O2	5.27	122.06	118.90
1	13	1197	G	OP1-P-O3'	5.27	116.79	105.20
26	1H	617	G	N3-C2-N2	5.27	123.59	119.90
26	1H	983	A	N7-C8-N9	-5.27	111.17	113.80
26	1H	1380	G	N1-C2-N3	5.27	127.06	123.90
26	1H	1633	G	OP2-P-O3'	5.27	116.79	105.20
26	1H	2561	A	N1-C2-N3	5.27	131.94	129.30
34	58	15	LEU	CA-CB-CG	5.27	127.42	115.30
1	13	413	G	C8-N9-C4	-5.27	104.29	106.40
26	1H	131	G	N1-C6-O6	5.27	123.06	119.90
26	1H	585	G	C5-C6-O6	-5.27	125.44	128.60
26	1H	612	G	N1-C6-O6	-5.27	116.74	119.90
26	1H	1021	A	C4-C5-N7	5.27	113.33	110.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	1899	G	C5-C6-O6	5.27	131.76	128.60
1	1G	1270	C	C5-C6-N1	5.27	123.63	121.00
1	1G	1374	A	O4'-C1'-N9	5.27	112.41	108.20
26	14	141(A)	C	C6-N1-C2	-5.27	118.19	120.30
26	14	773	U	C5-C6-N1	-5.27	120.07	122.70
26	14	1616	A	C6-C5-N7	-5.27	128.61	132.30
26	14	2329	G	C6-N1-C2	-5.27	121.94	125.10
1	13	550	G	N1-C2-N2	5.27	120.94	116.20
23	2K	76	C	N3-C4-N4	5.27	121.69	118.00
26	1H	837	C	C6-N1-C2	-5.27	118.19	120.30
1	1G	784	C	C6-N1-C2	5.27	122.41	120.30
26	14	1806	C	OP2-P-O3'	5.27	116.79	105.20
1	13	529	G	C4-C5-N7	5.26	112.91	110.80
26	1H	1394	U	O5'-P-OP2	5.26	117.02	110.70
26	1H	1926	U	O5'-P-OP2	-5.26	100.96	105.70
26	1H	2210	G	N7-C8-N9	5.26	115.73	113.10
26	1H	2866	U	C5-C4-O4	5.26	129.06	125.90
46	H8	24	LEU	CA-CB-CG	5.26	127.41	115.30
26	14	821	A	C6-C5-N7	-5.26	128.62	132.30
1	13	1099	G	C8-N9-C4	-5.26	104.30	106.40
26	1H	188	G	O5'-P-OP2	-5.26	100.96	105.70
26	1H	736	C	O5'-P-OP2	5.26	117.01	110.70
26	1H	1188	U	OP2-P-O3'	5.26	116.78	105.20
26	1H	1990	C	N3-C2-O2	-5.26	118.22	121.90
26	1H	2069	G	C8-N9-C4	5.26	108.50	106.40
26	1H	2830	G	N7-C8-N9	5.26	115.73	113.10
48	J8	41	ARG	NE-CZ-NH1	5.26	122.93	120.30
26	14	252	G	C4-C5-N7	-5.26	108.69	110.80
26	14	775	G	C6-N1-C2	-5.26	121.94	125.10
26	14	2542	A	C5-N7-C8	5.26	106.53	103.90
26	1H	621	A	C5-C6-N1	-5.26	115.07	117.70
26	1H	1819	A	N1-C6-N6	5.26	121.75	118.60
1	1G	388	G	C5-C6-O6	5.26	131.75	128.60
1	13	185	A	C8-N9-C4	-5.26	103.70	105.80
1	13	285	G	C8-N9-C4	5.26	108.50	106.40
26	14	756	C	C2-N3-C4	5.26	122.53	119.90
26	14	1556	C	C6-N1-C2	5.26	122.40	120.30
1	1G	1397	C	C5-C6-N1	5.25	123.63	121.00
26	14	415	A	O5'-P-OP2	-5.25	100.97	105.70
26	14	2523	G	N1-C6-O6	5.25	123.05	119.90
24	3K	71	G	C8-N9-C1'	5.25	133.83	127.00
26	1H	191	A	O5'-P-OP1	5.25	117.00	110.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	729	G	N3-C2-N2	-5.25	116.22	119.90
26	1H	911	A	C5-N7-C8	-5.25	101.27	103.90
26	1H	1408	C	OP1-P-O3'	5.25	116.76	105.20
26	1H	1566	A	O5'-P-OP1	5.25	117.00	110.70
26	1H	2424	C	C5-C6-N1	5.25	123.63	121.00
26	1H	2547	U	N3-C4-O4	5.25	123.08	119.40
26	1H	266	G	O5'-P-OP2	-5.25	100.97	105.70
26	1H	864	G	C2-N3-C4	5.25	114.53	111.90
26	1H	1357	U	C4-C5-C6	5.25	122.85	119.70
26	1H	2396	G	N3-C4-N9	-5.25	122.85	126.00
1	1G	28	G	N7-C8-N9	5.25	115.73	113.10
1	1G	1465	C	N1-C2-O2	5.25	122.05	118.90
26	14	476	G	N1-C6-O6	5.25	123.05	119.90
26	14	773	U	N1-C2-N3	5.25	118.05	114.90
26	14	2490	G	C8-N9-C4	-5.25	104.30	106.40
26	1H	782	A	C2-N3-C4	-5.25	107.97	110.60
1	1G	1487	G	N3-C2-N2	-5.25	116.22	119.90
26	14	2501	C	N3-C2-O2	5.25	125.58	121.90
26	14	2755	C	C2-N1-C1'	5.25	124.58	118.80
1	13	994	A	C2-N3-C4	5.25	113.22	110.60
26	1H	59	U	N3-C4-C5	-5.25	111.45	114.60
26	1H	791	C	P-O3'-C3'	5.25	126.00	119.70
26	1H	870	A	C5-C6-N1	5.25	120.32	117.70
26	1H	2056	G	OP1-P-O3'	5.25	116.75	105.20
26	1H	2491	U	N1-C2-N3	-5.25	111.75	114.90
26	14	97	C	O5'-P-OP2	-5.25	100.98	105.70
26	14	863	A	C8-N9-C4	-5.25	103.70	105.80
26	14	978	G	O5'-P-OP2	-5.25	100.98	105.70
26	14	1770	G	C8-N9-C4	-5.25	104.30	106.40
26	14	2038	G	C4-C5-N7	5.25	112.90	110.80
1	13	738	C	C5-C6-N1	5.25	123.62	121.00
26	1H	321	G	C4-N9-C1'	5.25	133.32	126.50
26	1H	2469	A	C4-C5-N7	5.25	113.32	110.70
1	1G	690	G	C4-C5-N7	5.25	112.90	110.80
23	2L	24	C	C6-N1-C2	-5.25	118.20	120.30
26	14	774	A	C5-C6-N1	-5.25	115.08	117.70
26	14	911	A	OP2-P-O3'	-5.25	93.66	105.20
1	13	1404	C	N1-C2-O2	-5.25	115.75	118.90
26	1H	624	C	C5-C4-N4	-5.25	116.53	120.20
26	1H	946	G	C5-N7-C8	5.25	106.92	104.30
26	14	2406	U	O4'-C1'-N1	-5.25	104.00	108.20
26	1H	937	U	N3-C2-O2	5.24	125.87	122.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	945	A	C6-N1-C2	-5.24	115.45	118.60
26	1H	1246	A	N1-C2-N3	5.24	131.92	129.30
26	1H	1369	G	C5-N7-C8	5.24	106.92	104.30
26	1H	1381	G	C4-C5-N7	-5.24	108.70	110.80
26	1H	1939	U	C4-C5-C6	-5.24	116.55	119.70
26	1H	2414	G	N3-C2-N2	-5.24	116.23	119.90
26	1H	2433	A	C4-C5-C6	5.24	119.62	117.00
26	1H	2580	U	OP2-P-O3'	5.24	116.74	105.20
26	1H	2731	G	C4-C5-N7	5.24	112.90	110.80
26	14	1334	G	O5'-P-OP1	-5.24	100.98	105.70
1	1G	893	C	N1-C2-O2	5.24	122.05	118.90
26	14	2070	G	C2-N3-C4	-5.24	109.28	111.90
1	13	1508	G	C6-C5-N7	5.24	133.54	130.40
26	1H	869	G	N3-C2-N2	5.24	123.57	119.90
26	1H	1464	C	O5'-P-OP1	-5.24	100.98	105.70
27	16	79	C	C5-C4-N4	-5.24	116.53	120.20
1	1G	115	G	P-O3'-C3'	5.24	125.99	119.70
26	14	1955	U	C5-C4-O4	5.24	129.04	125.90
1	13	452	A	C8-N9-C4	5.24	107.89	105.80
1	13	513	C	C2-N3-C4	5.24	122.52	119.90
26	1H	314	A	C4-C5-C6	-5.24	114.38	117.00
26	1H	1007	C	N1-C2-O2	-5.24	115.76	118.90
26	1H	1163	G	C8-N9-C4	-5.24	104.30	106.40
26	1H	1264	G	N3-C4-C5	-5.24	125.98	128.60
26	1H	1278	A	N9-C4-C5	-5.24	103.70	105.80
26	1H	1788	C	N1-C2-O2	5.24	122.04	118.90
26	1H	2085	C	N3-C2-O2	5.24	125.57	121.90
26	1H	2271	G	C8-N9-C1'	-5.24	120.19	127.00
1	1G	904	C	C6-N1-C2	-5.24	118.20	120.30
26	14	1805	U	N1-C2-N3	5.24	118.04	114.90
1	13	63	C	C5-C6-N1	5.24	123.62	121.00
1	1G	525	C	C5-C6-N1	5.24	123.62	121.00
26	14	1297	C	OP1-P-O3'	5.24	116.72	105.20
26	14	2437	U	N3-C4-O4	-5.24	115.73	119.40
26	1H	116	C	OP2-P-O3'	5.24	116.72	105.20
26	1H	409	C	C6-N1-C2	5.24	122.39	120.30
26	1H	831	G	N7-C8-N9	-5.24	110.48	113.10
26	1H	944	G	C4-N9-C1'	5.24	133.31	126.50
26	14	270(U)	C	C5-C6-N1	5.24	123.62	121.00
26	14	603	A	O4'-C1'-N9	5.24	112.39	108.20
26	14	776	G	N1-C6-O6	5.24	123.04	119.90
1	13	101	A	N1-C6-N6	5.23	121.74	118.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	13	387	U	OP1-P-O3'	5.23	116.71	105.20
26	1H	575	A	C6-N1-C2	-5.23	115.46	118.60
26	1H	128	C	C5-C6-N1	-5.23	118.38	121.00
26	1H	963	U	OP1-P-OP2	-5.23	111.75	119.60
26	1H	1200	C	N3-C4-C5	-5.23	119.81	121.90
26	1H	1681	G	C4-C5-N7	5.23	112.89	110.80
26	1H	1838	C	C6-N1-C2	5.23	122.39	120.30
26	1H	2870	C	OP2-P-O3'	5.23	116.71	105.20
1	1G	316	G	OP1-P-O3'	5.23	116.71	105.20
26	14	2235	G	N3-C4-C5	-5.23	125.98	128.60
26	1H	914	C	C6-N1-C1'	5.23	127.08	120.80
26	1H	946	G	OP1-P-OP2	-5.23	111.75	119.60
26	1H	2359	C	O5'-P-OP1	-5.23	100.99	105.70
26	14	2826	A	N1-C6-N6	-5.23	115.46	118.60
26	1H	1836	C	C6-N1-C2	-5.23	118.21	120.30
26	1H	2583	G	N7-C8-N9	5.23	115.71	113.10
26	14	624	C	N1-C2-O2	-5.23	115.76	118.90
26	14	1302	A	N9-C4-C5	5.23	107.89	105.80
26	14	1819	A	P-O3'-C3'	5.23	125.97	119.70
23	2K	63	C	N1-C2-O2	5.23	122.03	118.90
26	1H	586	A	C2-N3-C4	5.23	113.21	110.60
26	1H	2378	A	C8-N9-C4	5.23	107.89	105.80
26	1H	2506	U	P-O3'-C3'	5.23	125.97	119.70
26	14	1618	A	N9-C4-C5	5.23	107.89	105.80
26	1H	779	U	N3-C2-O2	5.22	125.86	122.20
26	1H	1571	A	C6-N1-C2	-5.22	115.47	118.60
26	1H	1899	G	N7-C8-N9	5.22	115.71	113.10
26	1H	2708	G	C5-C6-O6	-5.22	125.47	128.60
26	14	1001	A	O5'-P-OP1	-5.22	101.00	105.70
26	14	1811	G	C5-C6-O6	-5.22	125.47	128.60
1	13	23	C	N3-C4-C5	-5.22	119.81	121.90
1	13	949	A	N7-C8-N9	-5.22	111.19	113.80
1	13	1402	C	C5-C4-N4	5.22	123.86	120.20
26	1H	534	U	OP1-P-OP2	5.22	127.43	119.60
41	C8	74	LEU	CA-CB-CG	5.22	127.31	115.30
1	1G	380	G	N9-C4-C5	5.22	107.49	105.40
1	1G	667	G	N1-C6-O6	5.22	123.03	119.90
1	1G	1496	C	O5'-P-OP2	-5.22	101.00	105.70
1	13	949	A	C8-N9-C4	5.22	107.89	105.80
26	1H	1808	U	N3-C4-O4	5.22	123.05	119.40
26	14	819	A	N7-C8-N9	5.22	116.41	113.80
26	14	2497	A	O5'-P-OP2	5.22	116.97	110.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	1805	U	OP2-P-O3'	5.22	116.68	105.20
27	16	56	G	C4-N9-C1'	5.22	133.29	126.50
26	14	67	U	N3-C4-C5	-5.22	111.47	114.60
26	14	470	A	N1-C6-N6	5.22	121.73	118.60
26	14	835	A	C5-C6-N1	5.22	120.31	117.70
27	1J	14	U	OP1-P-OP2	5.22	127.43	119.60
26	1H	2074	U	C5-C4-O4	5.22	129.03	125.90
26	1H	2079	U	N3-C4-O4	5.22	123.05	119.40
26	14	270(Q)	C	C5-C6-N1	5.22	123.61	121.00
26	14	1672	C	N3-C4-C5	-5.22	119.81	121.90
26	14	1779	U	C6-N1-C1'	-5.22	113.89	121.20
26	14	2274	A	C5-N7-C8	-5.22	101.29	103.90
1	13	1502	A	C5-C6-N6	-5.22	119.53	123.70
26	1H	321	G	C8-N9-C1'	-5.22	120.22	127.00
26	1H	985	C	OP1-P-OP2	-5.22	111.78	119.60
26	1H	1391	U	N3-C4-O4	5.22	123.05	119.40
26	1H	2502	G	C5-C6-N1	5.22	114.11	111.50
27	16	50	G	OP2-P-O3'	5.22	116.68	105.20
26	14	35	G	N3-C2-N2	-5.22	116.25	119.90
27	1J	116	G	N1-C6-O6	5.22	123.03	119.90
26	1H	763	G	C4-C5-N7	5.21	112.89	110.80
26	1H	938	G	N1-C6-O6	-5.21	116.77	119.90
26	1H	1644	C	C2-N1-C1'	5.21	124.54	118.80
26	1H	2068	U	O5'-P-OP1	-5.21	101.01	105.70
26	1H	2337	G	OP1-P-OP2	-5.21	111.78	119.60
1	1G	345	C	C2-N1-C1'	5.21	124.54	118.80
26	14	680	G	N1-C6-O6	-5.21	116.77	119.90
26	14	797	C	N3-C4-C5	-5.21	119.81	121.90
26	14	801	G	C6-C5-N7	5.21	133.53	130.40
26	14	1324	G	C5-C6-O6	-5.21	125.47	128.60
1	13	1286	A	C8-N9-C4	-5.21	103.72	105.80
26	14	1590	U	O5'-P-OP1	-5.21	101.01	105.70
26	14	1997	G	C6-N1-C2	-5.21	121.97	125.10
26	14	2689	U	P-O3'-C3'	5.21	125.96	119.70
27	1J	12	C	N3-C2-O2	-5.21	118.25	121.90
1	13	1483	A	C5-N7-C8	-5.21	101.29	103.90
26	1H	397	G	N3-C4-C5	5.21	131.21	128.60
26	1H	835	A	C8-N9-C4	-5.21	103.72	105.80
26	1H	2494	G	N3-C2-N2	-5.21	116.25	119.90
27	16	60	C	N1-C2-O2	5.21	122.03	118.90
1	1G	666	G	C6-C5-N7	-5.21	127.27	130.40
26	14	800	A	OP1-P-OP2	-5.21	111.78	119.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	921	G	N9-C4-C5	5.21	107.48	105.40
26	14	1778	U	N3-C2-O2	5.21	125.85	122.20
26	14	2602	A	N1-C6-N6	-5.21	115.47	118.60
33	69	77	LEU	CA-CB-CG	5.21	127.28	115.30
26	1H	450	G	C5-C6-N1	-5.21	108.89	111.50
26	14	1672	C	N1-C2-O2	-5.21	115.77	118.90
26	14	2423	U	C2-N1-C1'	-5.21	111.45	117.70
26	1H	324	A	C4-C5-C6	5.21	119.60	117.00
26	1H	411	G	N3-C2-N2	-5.21	116.25	119.90
26	1H	600	G	C2-N3-C4	-5.21	109.30	111.90
26	1H	1274	A	C4-C5-C6	5.21	119.60	117.00
26	1H	1624	G	N7-C8-N9	-5.21	110.50	113.10
26	1H	1660	C	C6-N1-C2	-5.21	118.22	120.30
26	1H	2862	G	OP1-P-O3'	5.21	116.66	105.20
26	14	949	C	OP2-P-O3'	5.21	116.66	105.20
1	13	1358	U	C5-C4-O4	-5.21	122.78	125.90
26	1H	1939	U	C5-C4-O4	-5.21	122.78	125.90
26	14	83	G	C8-N9-C4	-5.21	104.32	106.40
26	14	2243	U	OP1-P-O3'	-5.21	93.74	105.20
1	13	721	G	N3-C4-C5	-5.21	126.00	128.60
26	1H	928	G	C4-C5-N7	5.21	112.88	110.80
26	1H	2060	A	C4-N9-C1'	-5.21	116.93	126.30
26	14	1216	G	C4-C5-C6	5.21	121.92	118.80
26	14	1605	C	OP1-P-OP2	5.21	127.41	119.60
23	2K	58	A	OP1-P-OP2	-5.20	111.79	119.60
26	1H	815	C	O5'-P-OP2	-5.20	101.02	105.70
26	1H	2215	G	C8-N9-C4	5.20	108.48	106.40
26	14	571	A	N1-C6-N6	5.20	121.72	118.60
26	14	1809	A	OP1-P-OP2	-5.20	111.79	119.60
26	14	1950	G	N1-C6-O6	-5.20	116.78	119.90
48	F5	29	GLY	N-CA-C	5.20	126.11	113.10
26	14	386	G	C6-C5-N7	-5.20	127.28	130.40
26	14	2700	C	N1-C2-O2	5.20	122.02	118.90
26	1H	740	U	OP1-P-O3'	-5.20	93.76	105.20
26	1H	1229(A)	G	C2-N3-C4	-5.20	109.30	111.90
26	14	573	G	N9-C4-C5	5.20	107.48	105.40
26	14	738	G	O5'-P-OP2	-5.20	101.02	105.70
26	14	2239	G	OP2-P-O3'	5.20	116.64	105.20
1	13	402	G	O5'-P-OP1	5.20	116.94	110.70
26	1H	579	G	C6-N1-C2	5.20	128.22	125.10
26	1H	845	G	C8-N9-C1'	5.20	133.76	127.00
26	1H	2327	A	C2-N3-C4	5.20	113.20	110.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	2330	G	N9-C4-C5	-5.20	103.32	105.40
26	1H	2490	G	O4'-C1'-N9	5.20	112.36	108.20
26	1H	2636	U	O5'-P-OP1	-5.20	101.02	105.70
26	14	646	A	C8-N9-C4	-5.20	103.72	105.80
1	13	571	U	C5-C6-N1	5.20	125.30	122.70
26	1H	729	G	C5-N7-C8	-5.20	101.70	104.30
26	1H	754	C	C2-N3-C4	-5.20	117.30	119.90
26	1H	950	G	N1-C6-O6	5.20	123.02	119.90
26	1H	953	A	O5'-P-OP2	-5.20	101.02	105.70
26	1H	2416	C	O5'-P-OP2	-5.20	101.02	105.70
26	14	2060	A	O4'-C1'-N9	5.20	112.36	108.20
26	1H	1936	A	C6-C5-N7	-5.20	128.66	132.30
26	1H	2072	G	OP1-P-OP2	-5.20	111.81	119.60
26	1H	2251	G	N3-C4-C5	-5.20	126.00	128.60
27	16	110	G	C8-N9-C4	-5.20	104.32	106.40
26	14	226	G	C6-C5-N7	-5.20	127.28	130.40
26	14	2544	G	N1-C6-O6	5.20	123.02	119.90
26	14	2607	G	N3-C4-N9	5.20	129.12	126.00
1	13	56	U	C6-N1-C2	-5.19	117.88	121.00
26	1H	807	U	OP1-P-OP2	5.19	127.39	119.60
1	13	893	C	C5-C4-N4	-5.19	116.56	120.20
26	1H	1613	G	OP2-P-O3'	5.19	116.62	105.20
26	1H	1673	U	C5-C6-N1	-5.19	120.10	122.70
26	1H	2034	U	C6-N1-C2	-5.19	117.88	121.00
26	14	133	C	N3-C4-C5	5.19	123.98	121.90
26	14	566	U	N3-C4-C5	5.19	117.72	114.60
26	14	2346	A	N9-C1'-C2'	5.19	120.75	114.00
26	1H	428	A	OP1-P-O3'	5.19	116.62	105.20
26	1H	659	C	C2-N1-C1'	-5.19	113.09	118.80
26	1H	2057	A	N1-C2-N3	5.19	131.90	129.30
26	1H	2415	G	C5-C6-O6	-5.19	125.48	128.60
26	14	677	A	O5'-P-OP1	-5.19	101.03	105.70
26	14	1773	A	OP1-P-O3'	5.19	116.62	105.20
26	1H	2373	G	N3-C4-N9	5.19	129.11	126.00
1	1G	1398	A	O4'-C1'-N9	5.19	112.35	108.20
26	14	2258	C	O5'-P-OP1	-5.19	101.03	105.70
26	1H	1195	G	C5-C6-O6	-5.19	125.49	128.60
26	14	511	U	N3-C2-O2	-5.19	118.57	122.20
26	14	2320	A	P-O3'-C3'	5.19	125.92	119.70
39	65	58	LEU	CA-CB-CG	5.19	127.23	115.30
1	13	550	G	N3-C2-N2	-5.19	116.27	119.90
1	13	576	G	C8-N9-C1'	-5.19	120.26	127.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	13	991	U	C2-N1-C1'	5.19	123.92	117.70
26	1H	404	C	P-O3'-C3'	5.19	125.92	119.70
26	1H	1900	A	C5'-C4'-O4'	-5.19	102.88	109.10
26	1H	631	A	N7-C8-N9	-5.18	111.21	113.80
26	1H	1303	G	N3-C2-N2	5.18	123.53	119.90
26	1H	1770	G	C8-N9-C4	5.18	108.47	106.40
26	1H	2485	G	N1-C2-N2	-5.18	111.53	116.20
26	14	399	G	N1-C2-N2	-5.18	111.53	116.20
54	L5	8	ASN	CB-CA-C	5.18	120.77	110.40
26	1H	2291	U	OP2-P-O3'	5.18	116.60	105.20
26	1H	2745	C	OP1-P-OP2	-5.18	111.83	119.60
26	1H	2822	G	N9-C4-C5	-5.18	103.33	105.40
1	1G	1489	G	O5'-P-OP2	-5.18	101.04	105.70
26	14	1258	C	OP2-P-O3'	5.18	116.60	105.20
26	14	1807	G	N1-C6-O6	5.18	123.01	119.90
26	14	1967	C	O5'-P-OP2	-5.18	101.04	105.70
26	14	2325	G	C8-N9-C4	-5.18	104.33	106.40
26	14	2331	G	O5'-P-OP2	-5.18	101.04	105.70
26	1H	2577	A	C5-C6-N6	5.18	127.84	123.70
26	1H	2709	G	O5'-P-OP1	5.18	116.92	110.70
26	14	1259	G	OP2-P-O3'	5.18	116.60	105.20
1	13	580	U	N3-C4-O4	-5.18	115.77	119.40
26	1H	77	C	C5-C4-N4	-5.18	116.57	120.20
26	1H	130	C	N3-C4-N4	5.18	121.63	118.00
26	1H	211	A	N9-C4-C5	-5.18	103.73	105.80
26	1H	629	G	O5'-P-OP1	5.18	116.92	110.70
26	1H	2294	C	OP1-P-OP2	-5.18	111.83	119.60
26	14	2380	C	C2-N3-C4	-5.18	117.31	119.90
26	14	2568	C	N3-C2-O2	-5.18	118.27	121.90
26	14	1337	G	OP1-P-O3'	5.18	116.59	105.20
1	13	1322	C	O5'-P-OP2	-5.18	101.04	105.70
1	13	1328	C	N3-C2-O2	5.18	125.52	121.90
26	1H	609	A	N9-C4-C5	-5.18	103.73	105.80
26	1H	2573	C	C6-N1-C2	-5.18	118.23	120.30
26	14	1969	A	OP1-P-OP2	-5.18	111.83	119.60
26	14	1982	C	C2-N3-C4	5.18	122.49	119.90
26	14	2001	A	C6-C5-N7	-5.18	128.68	132.30
26	14	2323	G	C8-N9-C4	5.18	108.47	106.40
1	13	1099	G	N9-C4-C5	5.17	107.47	105.40
1	13	1305	G	C5-C6-N1	-5.17	108.91	111.50
26	1H	46	C	OP2-P-O3'	5.17	116.58	105.20
1	1G	1054	C	N1-C2-O2	5.17	122.00	118.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	138	G	N7-C8-N9	5.17	115.69	113.10
26	14	1272	A	N1-C2-N3	-5.17	126.71	129.30
1	13	1527	C	C6-N1-C2	-5.17	118.23	120.30
26	1H	1368	G	N3-C4-C5	-5.17	126.01	128.60
26	1H	1416	G	O4'-C1'-N9	5.17	112.34	108.20
26	1H	2252	G	N3-C4-C5	5.17	131.19	128.60
1	1G	553	A	C8-N9-C4	-5.17	103.73	105.80
26	14	2278	A	OP1-P-O3'	5.17	116.58	105.20
26	1H	114	U	N1-C2-O2	5.17	126.42	122.80
26	1H	2256	G	OP2-P-O3'	5.17	116.58	105.20
26	1H	2331	G	N3-C4-C5	5.17	131.19	128.60
1	1G	41	G	C8-N9-C4	5.17	108.47	106.40
26	14	1544	C	N1-C2-O2	5.17	122.00	118.90
1	13	67	C	C6-N1-C2	-5.17	118.23	120.30
1	13	437	U	C6-N1-C2	-5.17	117.90	121.00
1	13	1058	G	OP1-P-OP2	-5.17	111.85	119.60
26	1H	202	U	C4-C5-C6	-5.17	116.60	119.70
26	1H	1640	C	C2-N1-C1'	-5.17	113.11	118.80
26	1H	2705	A	C5-N7-C8	-5.17	101.32	103.90
1	1G	1506	U	O5'-P-OP1	5.17	116.90	110.70
26	1H	952	G	N1-C2-N3	-5.17	120.80	123.90
26	1H	1425	G	N1-C6-O6	-5.17	116.80	119.90
1	1G	1301	U	N3-C2-O2	-5.17	118.58	122.20
26	14	1771	C	OP1-P-O3'	5.17	116.57	105.20
26	14	2256	G	N3-C4-N9	5.17	129.10	126.00
26	1H	257	A	C8-N9-C4	-5.17	103.73	105.80
26	1H	2325	G	C8-N9-C4	-5.17	104.33	106.40
23	2L	30	G	N3-C4-N9	5.17	129.10	126.00
26	14	1678	G	N1-C6-O6	5.17	123.00	119.90
27	16	115	G	N3-C4-C5	-5.16	126.02	128.60
1	1G	774	G	C4-C5-N7	5.16	112.86	110.80
26	14	155	C	N1-C2-O2	5.16	122.00	118.90
26	14	642	G	C5-N7-C8	-5.16	101.72	104.30
26	14	840	C	C6-N1-C2	5.16	122.37	120.30
26	14	1914	C	C2-N1-C1'	5.16	124.48	118.80
26	14	2433	A	OP2-P-O3'	5.16	116.56	105.20
1	1G	793	U	C2-N1-C1'	-5.16	111.51	117.70
26	14	833	U	N3-C4-O4	5.16	123.01	119.40
1	13	1205	U	N1-C2-O2	-5.16	119.19	122.80
1	13	1513	A	N1-C6-N6	5.16	121.70	118.60
26	1H	97	C	OP1-P-OP2	5.16	127.34	119.60
26	1H	698	C	OP1-P-OP2	5.16	127.34	119.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	1597	A	O4'-C1'-N9	5.16	112.33	108.20
26	14	595	C	C5-C6-N1	5.16	123.58	121.00
26	14	2244	U	C2-N1-C1'	-5.16	111.51	117.70
26	14	2365	G	C4-C5-N7	5.16	112.86	110.80
26	14	2423	U	C6-N1-C2	5.16	124.10	121.00
1	13	813	U	N3-C4-C5	5.16	117.69	114.60
1	13	1451	A	C8-N9-C4	-5.16	103.74	105.80
26	1H	755	C	N1-C2-O2	-5.16	115.80	118.90
26	14	828	U	C6-N1-C2	-5.16	117.91	121.00
26	14	1479	G	N1-C6-O6	5.16	123.00	119.90
26	1H	1670	C	C4-C5-C6	5.16	119.98	117.40
26	14	1372	U	N1-C2-N3	5.16	117.99	114.90
26	14	1970	A	C5-C6-N6	5.16	127.83	123.70
26	14	2404	C	O5'-P-OP1	-5.16	101.06	105.70
1	13	782	A	OP2-P-O3'	5.16	116.54	105.20
26	1H	564	C	OP1-P-O3'	5.16	116.54	105.20
26	1H	1674	G	C6-C5-N7	-5.16	127.31	130.40
26	1H	2030	A	N1-C6-N6	5.16	121.69	118.60
26	1H	2301	C	O5'-P-OP1	-5.16	101.06	105.70
26	1H	2548	G	C6-N1-C2	-5.16	122.01	125.10
26	14	828	U	N3-C4-O4	-5.16	115.79	119.40
26	14	2520	C	N1-C2-O2	-5.16	115.81	118.90
26	1H	982	C	C6-N1-C2	-5.15	118.24	120.30
26	1H	1144	G	OP1-P-O3'	5.15	116.54	105.20
26	14	1979	C	C6-N1-C2	-5.15	118.24	120.30
1	13	353	A	OP2-P-O3'	5.15	116.53	105.20
26	1H	302	C	C6-N1-C2	-5.15	118.24	120.30
26	1H	1252	G	O4'-C1'-N9	-5.15	104.08	108.20
26	1H	1891	G	N3-C2-N2	-5.15	116.29	119.90
26	14	2586	C	C5-C4-N4	-5.15	116.59	120.20
26	14	2699	C	N3-C2-O2	5.15	125.51	121.90
1	13	856	C	C5-C6-N1	5.15	123.58	121.00
26	1H	121	G	C6-C5-N7	-5.15	127.31	130.40
26	1H	751	A	C6-N1-C2	-5.15	115.51	118.60
26	1H	1417	C	C6-N1-C2	-5.15	118.24	120.30
26	1H	2089	U	N3-C4-O4	5.15	123.01	119.40
26	1H	2346	A	C5-C6-N1	-5.15	115.12	117.70
26	1H	2602	A	C2-N3-C4	5.15	113.17	110.60
26	1H	122	G	C5-C6-O6	-5.15	125.51	128.60
26	1H	132	G	C4-N9-C1'	5.15	133.19	126.50
26	1H	1982	C	C2-N1-C1'	5.15	124.46	118.80
26	14	1648	C	C6-N1-C2	-5.15	118.24	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	13	721	G	N3-C4-N9	5.15	129.09	126.00
26	1H	148	C	OP1-P-OP2	5.15	127.32	119.60
26	1H	432	A	C5-N7-C8	-5.15	101.33	103.90
26	1H	2665	A	N1-C2-N3	5.15	131.87	129.30
43	E8	64	MET	N-CA-C	5.15	124.90	111.00
1	1G	322	C	N3-C4-N4	5.15	121.60	118.00
26	14	1359	A	C4-C5-C6	-5.15	114.43	117.00
26	14	1695	G	N9-C4-C5	-5.15	103.34	105.40
26	1H	814	C	N3-C4-C5	5.15	123.96	121.90
26	1H	939	G	OP2-P-O3'	5.15	116.52	105.20
1	13	812	C	O5'-P-OP2	5.14	116.87	110.70
26	1H	1031	G	N1-C2-N2	-5.14	111.57	116.20
23	2L	41	C	C6-N1-C2	-5.14	118.24	120.30
26	14	459	U	C2-N1-C1'	5.14	123.87	117.70
26	14	518	G	N3-C2-N2	-5.14	116.30	119.90
26	14	1256	G	C6-C5-N7	-5.14	127.31	130.40
26	14	1644	C	N3-C2-O2	-5.14	118.30	121.90
26	14	1950	G	N3-C2-N2	5.14	123.50	119.90
26	1H	70	G	C5-C6-O6	5.14	131.69	128.60
1	1G	180	U	C5-C6-N1	5.14	125.27	122.70
1	1G	1441	G	O5'-P-OP1	-5.14	101.07	105.70
26	14	479	A	P-O3'-C3'	5.14	125.87	119.70
26	14	861	A	O5'-P-OP1	-5.14	101.07	105.70
26	14	1336	A	C6-N1-C2	-5.14	115.52	118.60
1	13	696	A	C8-N9-C4	-5.14	103.74	105.80
1	13	1406	U	OP2-P-O3'	5.14	116.51	105.20
26	14	929	G	C4-N9-C1'	5.14	133.18	126.50
26	14	1236	G	C8-N9-C4	5.14	108.46	106.40
26	14	1849	G	C8-N9-C4	-5.14	104.34	106.40
26	14	1930	G	C8-N9-C1'	5.14	133.68	127.00
26	1H	500	G	C4-C5-N7	-5.14	108.74	110.80
26	1H	748	G	C4-N9-C1'	-5.14	119.82	126.50
26	1H	840	C	C2-N1-C1'	-5.14	113.15	118.80
26	1H	2031	A	C5-C6-N6	-5.14	119.59	123.70
26	1H	2685	G	C4-C5-N7	-5.14	108.75	110.80
26	14	463	G	OP1-P-O3'	5.14	116.51	105.20
26	14	561	G	N3-C4-C5	5.14	131.17	128.60
26	14	621	A	N7-C8-N9	5.14	116.37	113.80
26	14	2584	U	O4'-C1'-N1	5.14	112.31	108.20
26	1H	321	G	C4-C5-C6	5.14	121.88	118.80
26	1H	543	C	C6-N1-C2	5.14	122.36	120.30
26	1H	1225	C	OP1-P-OP2	5.14	127.31	119.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	2495	G	C5-C6-O6	5.14	131.68	128.60
26	14	2508	G	C5-C6-N1	5.14	114.07	111.50
1	13	1518	A	C5-C6-N1	-5.14	115.13	117.70
26	1H	30	G	N1-C6-O6	-5.14	116.82	119.90
26	1H	2588	G	OP1-P-OP2	-5.14	111.90	119.60
26	14	391	G	C6-C5-N7	-5.14	127.32	130.40
26	14	982	C	C5-C6-N1	5.14	123.57	121.00
26	14	1728	G	N3-C4-N9	5.14	129.08	126.00
23	2K	9	G	C2-N3-C4	5.13	114.47	111.90
26	1H	719	C	C6-N1-C2	-5.13	118.25	120.30
26	1H	1643	G	OP2-P-O3'	5.13	116.50	105.20
26	1H	2713	A	N3-C4-C5	5.13	130.39	126.80
1	1G	769	G	C4-N9-C1'	5.13	133.17	126.50
15	6A	23	GLY	N-CA-C	5.13	125.94	113.10
26	14	561	G	N9-C4-C5	5.13	107.45	105.40
1	13	1332	A	N7-C8-N9	5.13	116.37	113.80
23	2K	31	G	C5-C6-N1	-5.13	108.93	111.50
26	1H	640	C	N3-C4-C5	-5.13	119.85	121.90
26	1H	1625	C	N1-C2-O2	5.13	121.98	118.90
26	1H	2364	C	OP2-P-O3'	5.13	116.49	105.20
26	14	2374	C	C5-C6-N1	-5.13	118.43	121.00
26	1H	655	A	N7-C8-N9	5.13	116.37	113.80
26	1H	919	G	C4-C5-C6	5.13	121.88	118.80
26	1H	1579	A	O5'-P-OP2	-5.13	101.08	105.70
26	14	1409	C	OP1-P-OP2	5.13	127.30	119.60
26	14	1899	G	P-O3'-C3'	5.13	125.86	119.70
26	14	1970	A	N1-C6-N6	-5.13	115.52	118.60
26	14	2673	G	C4-N9-C1'	5.13	133.17	126.50
26	1H	965	C	OP1-P-OP2	5.13	127.30	119.60
26	1H	1781	C	N3-C4-N4	-5.13	114.41	118.00
26	1H	792	G	C6-C5-N7	-5.13	127.32	130.40
26	1H	814	C	C6-N1-C2	5.13	122.35	120.30
26	1H	865	C	OP1-P-OP2	-5.13	111.91	119.60
26	1H	1284	A	N1-C2-N3	5.13	131.86	129.30
1	1G	650	G	N3-C2-N2	-5.13	116.31	119.90
26	14	791	C	C5-C4-N4	-5.13	116.61	120.20
26	14	1301	A	O4'-C1'-N9	5.13	112.30	108.20
1	13	771	G	C5-C6-N1	-5.13	108.94	111.50
1	13	1290	G	N1-C6-O6	5.13	122.97	119.90
26	1H	471	A	C4-C5-N7	5.13	113.26	110.70
26	1H	1281	G	C5-C6-O6	-5.13	125.53	128.60
26	1H	1297	C	N3-C2-O2	-5.13	118.31	121.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	1377	G	C4-C5-C6	5.13	121.88	118.80
26	1H	2079	U	N3-C4-C5	-5.13	111.52	114.60
26	1H	2373	G	N1-C2-N3	5.13	126.98	123.90
26	14	441	U	O5'-P-OP1	-5.13	101.09	105.70
26	14	528	A	N3-C4-C5	5.13	130.39	126.80
26	14	733	G	OP1-P-OP2	5.13	127.29	119.60
26	14	947	G	C5-C6-N1	-5.13	108.94	111.50
26	14	2321	G	C4-N9-C1'	5.13	133.16	126.50
26	14	2574	G	C8-N9-C4	-5.13	104.35	106.40
1	13	904	C	N3-C2-O2	-5.12	118.31	121.90
27	16	110	G	N9-C4-C5	5.12	107.45	105.40
26	14	2700	C	C2-N1-C1'	5.12	124.44	118.80
1	13	235	C	N3-C4-N4	5.12	121.59	118.00
1	13	783	C	C5-C6-N1	-5.12	118.44	121.00
1	13	1501	C	N3-C4-C5	-5.12	119.85	121.90
26	1H	826	U	N3-C4-C5	-5.12	111.53	114.60
26	1H	942	G	N1-C2-N2	5.12	120.81	116.20
26	1H	1816	G	C4-N9-C1'	-5.12	119.84	126.50
26	1H	2048	G	C5-C6-N1	-5.12	108.94	111.50
26	14	1728	G	C2-N3-C4	5.12	114.46	111.90
26	14	2504	U	OP1-P-OP2	-5.12	111.92	119.60
27	1J	61	G	N1-C6-O6	5.12	122.97	119.90
26	1H	293	U	N3-C4-O4	5.12	122.98	119.40
26	1H	1667	G	O5'-P-OP1	-5.12	101.09	105.70
1	1G	1433	A	O5'-P-OP1	-5.12	101.09	105.70
26	14	725	G	N3-C2-N2	-5.12	116.31	119.90
26	14	2587	A	C5-C6-N6	-5.12	119.60	123.70
26	1H	1768	U	OP2-P-O3'	5.12	116.46	105.20
26	1H	2713	A	OP2-P-O3'	5.12	116.47	105.20
26	14	479	A	N1-C6-N6	-5.12	115.53	118.60
1	13	129	U	C2-N1-C1'	-5.12	111.56	117.70
26	1H	2048	G	N7-C8-N9	5.12	115.66	113.10
26	1H	2269	A	N1-C6-N6	5.12	121.67	118.60
26	1H	2351	G	C6-C5-N7	-5.12	127.33	130.40
1	1G	1274	G	N7-C8-N9	5.12	115.66	113.10
26	14	1022	G	P-O3'-C3'	5.12	125.84	119.70
26	14	2332	U	C5-C6-N1	-5.12	120.14	122.70
41	85	20	LEU	CA-CB-CG	5.12	127.07	115.30
26	14	388	G	C5-C6-O6	5.12	131.67	128.60
26	14	593	G	N7-C8-N9	-5.12	110.54	113.10
26	14	2518	A	N9-C4-C5	-5.12	103.75	105.80
1	13	169	C	C6-N1-C2	-5.12	118.25	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	13	895	G	O5'-P-OP2	-5.12	101.10	105.70
26	1H	868	U	C4-C5-C6	5.12	122.77	119.70
26	1H	1012	U	N3-C2-O2	5.12	125.78	122.20
26	1H	1124	C	C5-C6-N1	-5.12	118.44	121.00
26	1H	1660	C	N1-C2-O2	5.12	121.97	118.90
26	1H	2335	A	O5'-P-OP1	-5.12	101.10	105.70
26	14	2352	A	N1-C2-N3	5.12	131.86	129.30
23	2K	11	A	OP2-P-O3'	5.11	116.45	105.20
26	1H	204	A	C6-N1-C2	-5.11	115.53	118.60
26	1H	2324	C	O5'-P-OP2	-5.11	101.10	105.70
26	1H	2456	C	C4-C5-C6	5.11	119.96	117.40
25	4L	21	C	C2-N1-C1'	5.11	124.42	118.80
26	14	502	A	N1-C2-N3	5.11	131.86	129.30
26	14	677	A	C8-N9-C4	-5.11	103.75	105.80
26	14	775	G	N1-C2-N2	-5.11	111.60	116.20
26	1H	2040	C	N3-C2-O2	5.11	125.48	121.90
27	16	61	G	N7-C8-N9	5.11	115.66	113.10
26	14	683	C	C2-N3-C4	-5.11	117.34	119.90
1	13	975	A	N1-C6-N6	5.11	121.67	118.60
26	1H	1403	C	N3-C2-O2	-5.11	118.32	121.90
26	14	784	A	C5-C6-N6	5.11	127.79	123.70
26	14	1280	G	OP2-P-O3'	5.11	116.44	105.20
27	1J	103	U	C6-N1-C2	5.11	124.07	121.00
1	13	1195	C	N1-C2-O2	-5.11	115.83	118.90
26	1H	451	C	N1-C2-O2	-5.11	115.83	118.90
26	1H	1357	U	N1-C2-N3	5.11	117.97	114.90
26	1H	1767	C	O5'-P-OP1	-5.11	101.10	105.70
1	1G	380	G	C8-N9-C4	-5.11	104.36	106.40
1	13	311	C	C6-N1-C2	-5.11	118.26	120.30
1	13	1331	G	C8-N9-C4	-5.11	104.36	106.40
26	1H	467	G	N1-C6-O6	5.11	122.96	119.90
26	1H	557	U	O5'-P-OP2	-5.11	101.10	105.70
26	1H	826	U	C5-C6-N1	-5.11	120.15	122.70
26	1H	974(A)	C	C2-N1-C1'	5.11	124.42	118.80
26	1H	1688	U	O5'-P-OP2	-5.11	101.10	105.70
26	1H	1910	G	O5'-P-OP1	-5.11	101.10	105.70
1	1G	831	U	C6-N1-C2	-5.11	117.94	121.00
26	14	751	A	OP1-P-OP2	-5.11	111.94	119.60
26	14	1726	G	C8-N9-C4	-5.11	104.36	106.40
1	13	1213	A	OP2-P-O3'	5.11	116.43	105.20
22	1K	34	G	C8-N9-C4	-5.11	104.36	106.40
26	1H	534	U	OP2-P-O3'	5.11	116.43	105.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	563	G	C8-N9-C4	5.11	108.44	106.40
27	16	73	A	N1-C6-N6	-5.11	115.54	118.60
23	2K	11	A	N1-C6-N6	-5.10	115.54	118.60
26	14	1816	G	O5'-P-OP2	5.10	116.83	110.70
22	1K	33	U	N3-C4-C5	-5.10	111.54	114.60
23	2K	11	A	N9-C4-C5	5.10	107.84	105.80
1	1G	924	C	OP1-P-OP2	5.10	127.25	119.60
26	1H	126	A	OP1-P-O3'	-5.10	93.98	105.20
26	1H	1192	G	N9-C4-C5	-5.10	103.36	105.40
26	1H	2199	A	OP2-P-O3'	5.10	116.42	105.20
26	14	1625	C	O5'-P-OP2	-5.10	101.11	105.70
26	1H	705	A	N9-C4-C5	-5.10	103.76	105.80
26	1H	1357	U	C6-N1-C2	-5.10	117.94	121.00
26	1H	2595	G	C5-N7-C8	-5.10	101.75	104.30
1	1G	1500	A	N1-C6-N6	5.10	121.66	118.60
26	14	621	A	C6-N1-C2	5.10	121.66	118.60
1	13	50	A	P-O3'-C3'	5.10	125.82	119.70
26	1H	199	A	C2-N3-C4	5.10	113.15	110.60
26	1H	237	C	N3-C4-C5	-5.10	119.86	121.90
26	1H	330	A	C4-C5-N7	5.10	113.25	110.70
26	1H	709	U	N1-C2-O2	-5.10	119.23	122.80
26	1H	915	C	OP1-P-OP2	-5.10	111.95	119.60
26	1H	978	G	OP1-P-O3'	5.10	116.42	105.20
26	1H	2075	U	N1-C2-N3	5.10	117.96	114.90
26	14	921	G	N3-C2-N2	-5.10	116.33	119.90
1	13	1479	C	C5-C4-N4	-5.10	116.63	120.20
26	1H	1310	G	N1-C6-O6	5.10	122.96	119.90
26	14	2887	U	O5'-P-OP1	-5.10	101.11	105.70
1	13	23	C	C2-N3-C4	5.09	122.45	119.90
1	13	302	G	N3-C4-C5	-5.09	126.05	128.60
23	2K	25	U	N3-C2-O2	-5.09	118.63	122.20
26	1H	526	A	C5-C6-N6	5.09	127.78	123.70
26	1H	631	A	OP1-P-O3'	5.09	116.41	105.20
26	1H	1142(A)	A	C5-N7-C8	-5.09	101.35	103.90
26	1H	1594	G	N3-C2-N2	-5.09	116.33	119.90
26	1H	2099	U	C6-N1-C2	-5.09	117.94	121.00
26	1H	2430	A	N7-C8-N9	5.09	116.35	113.80
32	51	172	LYS	C-N-CA	5.09	143.39	122.00
26	14	792	G	OP2-P-O3'	5.09	116.41	105.20
26	14	2829	C	N1-C2-O2	-5.09	115.84	118.90
1	13	974	A	N9-C4-C5	-5.09	103.76	105.80
26	1H	1386	C	C5-C6-N1	5.09	123.55	121.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	1669	A	C5-N7-C8	-5.09	101.35	103.90
26	1H	2578	G	OP2-P-O3'	5.09	116.41	105.20
26	14	1267	U	OP2-P-O3'	5.09	116.41	105.20
1	13	330	C	C6-N1-C2	-5.09	118.26	120.30
1	13	449	C	C2-N1-C1'	5.09	124.40	118.80
1	13	1290	G	C6-C5-N7	-5.09	127.34	130.40
26	1H	1982	C	OP1-P-OP2	5.09	127.24	119.60
1	1G	314	C	C6-N1-C2	-5.09	118.26	120.30
1	1G	413	G	N1-C6-O6	-5.09	116.85	119.90
1	1G	1260	C	C5-C6-N1	5.09	123.55	121.00
26	14	113	G	C8-N9-C4	5.09	108.44	106.40
1	13	918	A	N1-C6-N6	5.09	121.65	118.60
26	1H	68	G	O5'-P-OP1	-5.09	101.12	105.70
26	1H	701	G	C6-C5-N7	-5.09	127.35	130.40
26	1H	1204	A	C3'-C2'-C1'	-5.09	97.43	101.50
26	1H	1814	G	C6-C5-N7	-5.09	127.35	130.40
26	1H	2609	U	N1-C2-N3	5.09	117.95	114.90
26	14	537	C	N1-C2-O2	5.09	121.95	118.90
26	14	1322	A	OP2-P-O3'	5.09	116.40	105.20
26	14	1339	G	C2-N3-C4	-5.09	109.36	111.90
26	14	2318	G	C8-N9-C4	-5.09	104.36	106.40
1	13	527	G	C8-N9-C4	-5.09	104.36	106.40
26	1H	2589	A	C8-N9-C4	5.09	107.83	105.80
26	14	1680	U	OP1-P-O3'	5.09	116.39	105.20
26	14	1681	G	C2-N3-C4	-5.09	109.36	111.90
26	1H	1252	G	C4-C5-N7	-5.09	108.77	110.80
26	14	1394	U	OP1-P-OP2	-5.09	111.97	119.60
26	14	1844	C	OP1-P-OP2	-5.09	111.97	119.60
26	1H	1344	G	C5-N7-C8	-5.08	101.76	104.30
26	1H	1858	G	C6-C5-N7	-5.08	127.35	130.40
26	1H	2011	U	C2-N1-C1'	-5.08	111.60	117.70
26	1H	2638	G	N3-C4-C5	-5.08	126.06	128.60
26	14	1261	C	C2-N3-C4	-5.08	117.36	119.90
26	14	2330	G	N1-C2-N2	-5.08	111.62	116.20
24	3K	76	A	O5'-P-OP1	-5.08	101.12	105.70
26	1H	828	U	N1-C2-N3	-5.08	111.85	114.90
26	1H	1428	C	C2-N3-C4	-5.08	117.36	119.90
26	1H	2487	G	C5-C6-O6	-5.08	125.55	128.60
23	2L	40	C	O5'-P-OP1	-5.08	101.12	105.70
26	14	141	A	O4'-C1'-N9	5.08	112.27	108.20
26	14	774	A	C4-C5-N7	5.08	113.24	110.70
26	14	915	C	N3-C2-O2	-5.08	118.34	121.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	1383	C	N3-C2-O2	5.08	125.46	121.90
26	14	1801	G	C5-C6-O6	-5.08	125.55	128.60
26	14	2715	C	O5'-P-OP1	5.08	116.80	110.70
26	1H	609	A	N7-C8-N9	-5.08	111.26	113.80
26	14	632	A	OP1-P-OP2	-5.08	111.98	119.60
26	14	792	G	N3-C4-N9	5.08	129.05	126.00
26	1H	695	G	N1-C6-O6	-5.08	116.85	119.90
26	1H	1156	A	O5'-P-OP2	-5.08	101.13	105.70
26	1H	1258	C	C6-N1-C1'	5.08	126.89	120.80
26	1H	1791	A	C2-N3-C4	5.08	113.14	110.60
26	1H	2710	C	O5'-P-OP2	5.08	116.80	110.70
26	14	581	C	C6-N1-C2	-5.08	118.27	120.30
1	13	628	G	N1-C6-O6	5.08	122.95	119.90
26	1H	2300	G	N3-C4-C5	-5.08	126.06	128.60
26	1H	2571	C	N3-C4-C5	-5.08	119.87	121.90
26	14	1391	U	C5-C4-O4	-5.08	122.85	125.90
1	13	535	A	C5-C6-N6	5.08	127.76	123.70
26	1H	1139	G	O5'-P-OP2	-5.08	101.13	105.70
26	1H	2197	U	OP2-P-O3'	5.08	116.37	105.20
26	1H	2827	C	N1-C2-O2	-5.08	115.85	118.90
26	14	2579	C	C5-C4-N4	-5.08	116.65	120.20
1	13	304	U	N3-C4-C5	-5.08	111.56	114.60
23	2K	73	A	C8-N9-C4	5.08	107.83	105.80
26	1H	330	A	C6-C5-N7	-5.08	128.75	132.30
26	14	1309	G	C8-N9-C4	5.08	108.43	106.40
1	13	413	G	N9-C4-C5	5.07	107.43	105.40
26	1H	115	C	N1-C2-O2	-5.07	115.86	118.90
26	1H	622	G	C5-N7-C8	5.07	106.84	104.30
26	1H	1792	G	OP1-P-O3'	5.07	116.36	105.20
26	1H	1800	C	C4-C5-C6	5.07	119.94	117.40
26	1H	2243	U	OP1-P-O3'	-5.07	94.04	105.20
26	1H	2379	G	N9-C4-C5	-5.07	103.37	105.40
27	16	45	A	OP1-P-OP2	5.07	127.21	119.60
1	1G	261	U	N3-C2-O2	-5.07	118.65	122.20
26	14	580	C	C6-N1-C2	-5.07	118.27	120.30
26	1H	593	G	N1-C2-N3	5.07	126.94	123.90
26	1H	988	A	OP2-P-O3'	5.07	116.36	105.20
26	1H	1733	G	C5-C6-N1	-5.07	108.96	111.50
26	1H	2587	A	C5-C6-N6	5.07	127.76	123.70
26	1H	2261	C	O5'-P-OP2	-5.07	101.14	105.70
1	1G	284	G	C5-C6-O6	-5.07	125.56	128.60
26	14	1822	G	C5-C6-N1	-5.07	108.96	111.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	13	320	C	N3-C2-O2	5.07	125.45	121.90
26	1H	59	U	C5-C4-O4	5.07	128.94	125.90
26	1H	692	C	N3-C2-O2	5.07	125.45	121.90
26	1H	738	G	N1-C2-N2	-5.07	111.64	116.20
1	1G	1431	C	N3-C4-N4	5.07	121.55	118.00
26	14	602	G	N9-C4-C5	-5.07	103.37	105.40
27	1J	54	G	N7-C8-N9	5.07	115.63	113.10
1	13	792	A	C1'-O4'-C4'	-5.07	105.84	109.90
1	13	1126	U	P-O3'-C3'	5.07	125.78	119.70
26	1H	682	G	N9-C4-C5	-5.07	103.37	105.40
26	1H	1671	U	N3-C4-O4	5.07	122.95	119.40
26	14	1664	A	N1-C6-N6	-5.07	115.56	118.60
26	14	2314	C	N1-C2-O2	5.07	121.94	118.90
26	14	2510	C	O5'-P-OP2	-5.07	101.14	105.70
26	1H	2323	G	N7-C8-N9	5.07	115.63	113.10
1	1G	115	G	C5-C6-O6	5.07	131.64	128.60
1	1G	1498	U	N3-C2-O2	-5.07	118.65	122.20
26	14	806	C	C5-C4-N4	-5.07	116.65	120.20
26	14	1314	C	C6-N1-C1'	-5.07	114.72	120.80
26	14	1327	C	N3-C4-C5	-5.07	119.87	121.90
26	14	2365	G	C6-C5-N7	-5.07	127.36	130.40
26	14	2840	C	N3-C4-C5	-5.07	119.87	121.90
26	14	2859	G	N3-C4-C5	-5.07	126.07	128.60
1	13	667	G	C8-N9-C4	-5.06	104.37	106.40
1	13	1224	G	OP1-P-OP2	-5.06	112.00	119.60
26	1H	290	G	N3-C4-C5	-5.06	126.07	128.60
26	1H	940	G	N3-C4-N9	5.06	129.04	126.00
1	1G	180	U	N3-C4-O4	5.06	122.94	119.40
26	14	1274	A	N3-C4-N9	5.06	131.45	127.40
1	13	805	C	OP1-P-OP2	-5.06	112.00	119.60
26	1H	665	C	N1-C2-O2	-5.06	115.86	118.90
26	1H	913	U	C4-C5-C6	-5.06	116.66	119.70
26	1H	1260	G	C5-N7-C8	5.06	106.83	104.30
26	1H	1362	C	N3-C2-O2	5.06	125.44	121.90
26	1H	2392	A	O5'-P-OP1	-5.06	101.14	105.70
27	16	103	U	OP2-P-O3'	5.06	116.34	105.20
26	14	1757	U	N3-C4-C5	-5.06	111.56	114.60
1	13	535	A	C8-N9-C4	-5.06	103.78	105.80
26	1H	809	G	N7-C8-N9	-5.06	110.57	113.10
26	1H	2271	G	C4-N9-C1'	5.06	133.08	126.50
26	14	2225	A	P-O3'-C3'	5.06	125.77	119.70
26	14	2419	U	OP1-P-O3'	5.06	116.33	105.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	13	328	C	O5'-P-OP1	-5.06	101.15	105.70
1	13	1335	C	C5-C6-N1	-5.06	118.47	121.00
1	13	1432	G	N3-C4-N9	5.06	129.03	126.00
26	1H	129	C	OP2-P-O3'	5.06	116.33	105.20
26	1H	243	U	O5'-P-OP2	-5.06	101.15	105.70
26	1H	938	G	C5-C6-O6	5.06	131.64	128.60
26	1H	1204	A	C4-C5-C6	5.06	119.53	117.00
26	1H	1431	U	C4-C5-C6	-5.06	116.67	119.70
26	1H	1625	C	O5'-P-OP1	5.06	116.77	110.70
26	1H	1692	U	C2-N1-C1'	5.06	123.77	117.70
26	1H	1817	G	N1-C6-O6	-5.06	116.86	119.90
26	1H	1964	G	OP2-P-O3'	5.06	116.33	105.20
26	1H	2300	G	C6-C5-N7	-5.06	127.36	130.40
26	1H	2388	A	O4'-C1'-N9	5.06	112.25	108.20
26	1H	2451	A	C5-C6-N6	5.06	127.75	123.70
1	1G	898	G	O5'-P-OP2	-5.06	101.15	105.70
26	14	511	U	N3-C4-C5	-5.06	111.56	114.60
26	14	2427	C	N3-C2-O2	5.06	125.44	121.90
26	14	2689	U	C5-C4-O4	5.06	128.94	125.90
1	13	191(F)	U	C5-C6-N1	5.06	125.23	122.70
26	1H	26	G	N3-C4-C5	-5.06	126.07	128.60
26	1H	536	A	N1-C6-N6	-5.06	115.56	118.60
26	1H	1300	U	C6-N1-C2	-5.06	117.97	121.00
26	1H	1792	G	C5-C6-O6	5.06	131.63	128.60
26	1H	2823	A	N1-C6-N6	5.06	121.63	118.60
1	1G	1259	C	C5-C6-N1	5.06	123.53	121.00
26	14	1274	A	N9-C4-C5	-5.06	103.78	105.80
26	14	2612	C	O5'-P-OP2	-5.06	101.15	105.70
26	14	2732	G	C8-N9-C4	-5.06	104.38	106.40
1	13	574	A	OP1-P-OP2	-5.06	112.02	119.60
26	1H	686	G	O5'-P-OP2	-5.06	101.15	105.70
26	1H	871	U	C2-N1-C1'	5.06	123.77	117.70
26	1H	1354	A	C2-N3-C4	-5.06	108.07	110.60
26	1H	2557	G	C8-N9-C4	-5.06	104.38	106.40
26	1H	651	G	N1-C6-O6	-5.05	116.87	119.90
26	1H	1034	G	C5-C6-O6	-5.05	125.57	128.60
26	1H	1936	A	C5-C6-N1	5.05	120.23	117.70
26	1H	1987	G	N3-C2-N2	-5.05	116.36	119.90
1	1G	1356	G	N7-C8-N9	5.05	115.63	113.10
26	14	201	C	C2-N1-C1'	-5.05	113.24	118.80
26	14	205	G	N3-C2-N2	5.05	123.44	119.90
26	14	1772	G	OP1-P-OP2	5.05	127.18	119.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	13	552	U	O5'-P-OP2	-5.05	101.15	105.70
26	1H	62	C	C4-C5-C6	5.05	119.93	117.40
26	1H	146	G	C6-C5-N7	-5.05	127.37	130.40
26	1H	579	G	OP2-P-O3'	5.05	116.32	105.20
26	1H	618(A)	C	C6-N1-C2	5.05	122.32	120.30
26	1H	1408	C	N1-C2-O2	-5.05	115.87	118.90
26	1H	1526	G	C6-C5-N7	-5.05	127.37	130.40
26	1H	1977	A	C2-N3-C4	-5.05	108.07	110.60
26	14	1227	A	N1-C6-N6	-5.05	115.57	118.60
26	14	2627	G	N3-C4-C5	-5.05	126.07	128.60
26	1H	483	A	C8-N9-C4	-5.05	103.78	105.80
26	1H	621	A	C8-N9-C4	-5.05	103.78	105.80
26	1H	1024	G	C4-N9-C1'	5.05	133.07	126.50
26	1H	1204	A	N7-C8-N9	5.05	116.33	113.80
26	1H	1826	G	OP1-P-O3'	5.05	116.31	105.20
26	1H	1888	G	C8-N9-C1'	-5.05	120.43	127.00
26	14	134	C	N3-C4-C5	5.05	123.92	121.90
26	14	503	A	N1-C6-N6	-5.05	115.57	118.60
26	14	1610	A	N9-C4-C5	-5.05	103.78	105.80
27	1J	14	U	O5'-P-OP2	-5.05	101.15	105.70
1	13	768	A	N1-C2-N3	5.05	131.82	129.30
26	1H	645	C	C6-N1-C2	-5.05	118.28	120.30
26	1H	907	U	C5-C6-N1	-5.05	120.17	122.70
26	1H	2246	G	N7-C8-N9	-5.05	110.58	113.10
26	1H	2280	G	OP1-P-O3'	5.05	116.31	105.20
27	16	5	C	N3-C4-C5	5.05	123.92	121.90
39	A8	110	LEU	CA-CB-CG	-5.05	103.69	115.30
26	14	1210	A	C2-N3-C4	-5.05	108.08	110.60
26	14	2370	G	C8-N9-C4	5.05	108.42	106.40
26	14	2736	G	N1-C6-O6	5.05	122.93	119.90
26	1H	208	C	OP1-P-OP2	5.05	127.17	119.60
26	1H	418	G	N9-C4-C5	-5.05	103.38	105.40
26	1H	446	G	C4-N9-C1'	5.05	133.06	126.50
26	1H	928	G	C5-N7-C8	-5.05	101.78	104.30
26	1H	1241	A	C4-C5-N7	5.05	113.22	110.70
1	13	749	C	N1-C2-O2	5.04	121.93	118.90
26	1H	828	U	C2-N3-C4	5.04	130.03	127.00
26	1H	1783	A	C5-C6-N1	-5.04	115.18	117.70
26	1H	2439	A	C5-N7-C8	-5.04	101.38	103.90
26	14	734	A	C4-C5-C6	5.04	119.52	117.00
26	14	1216	G	C5-C6-N1	-5.04	108.98	111.50
26	14	1527	G	N1-C2-N2	5.04	120.74	116.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	13	229	U	C5-C6-N1	5.04	125.22	122.70
26	1H	49	A	C8-N9-C4	5.04	107.82	105.80
26	1H	383	U	O4'-C1'-N1	5.04	112.23	108.20
26	1H	1933	G	C8-N9-C4	-5.04	104.38	106.40
1	1G	250	A	P-O3'-C3'	5.04	125.75	119.70
26	14	191	A	N1-C2-N3	-5.04	126.78	129.30
26	14	2038	G	N3-C2-N2	5.04	123.43	119.90
26	14	2356	C	N3-C2-O2	5.04	125.43	121.90
26	14	2433	A	C5-N7-C8	-5.04	101.38	103.90
26	14	2719	G	C5-C6-N1	-5.04	108.98	111.50
1	13	1058	G	C8-N9-C4	5.04	108.42	106.40
26	1H	1269	A	C5-N7-C8	-5.04	101.38	103.90
26	1H	1602	U	C4-C5-C6	5.04	122.72	119.70
26	1H	1773	A	C6-N1-C2	5.04	121.62	118.60
27	16	56	G	N3-C4-C5	-5.04	126.08	128.60
26	14	216	A	OP1-P-O3'	5.04	116.28	105.20
26	14	995	C	C6-N1-C1'	5.04	126.85	120.80
26	14	2033	A	C5-C6-N1	5.04	120.22	117.70
26	14	2281	C	C2-N1-C1'	5.04	124.34	118.80
1	13	912	C	O5'-P-OP1	-5.04	101.17	105.70
26	1H	1924	C	N3-C4-C5	5.04	123.92	121.90
26	1H	2702	U	N3-C2-O2	-5.04	118.67	122.20
26	1H	2782	G	N7-C8-N9	5.04	115.62	113.10
34	58	120	LEU	CA-CB-CG	5.04	126.89	115.30
26	14	1681	G	C6-C5-N7	-5.04	127.38	130.40
1	13	1249	C	C6-N1-C2	-5.04	118.29	120.30
1	13	1498	U	C6-N1-C1'	-5.04	114.15	121.20
26	1H	486	C	O5'-P-OP2	5.04	116.74	110.70
26	14	111	A	N1-C6-N6	5.04	121.62	118.60
26	14	209	C	C2-N1-C1'	5.04	124.34	118.80
1	13	1260	C	C5-C6-N1	5.03	123.52	121.00
26	1H	471	A	N1-C6-N6	5.03	121.62	118.60
26	1H	937	U	C5-C6-N1	-5.03	120.18	122.70
26	1H	2450	A	N1-C2-N3	5.03	131.82	129.30
55	Q8	34	TRP	N-CA-C	5.03	124.59	111.00
26	14	790	C	OP2-P-O3'	5.03	116.27	105.20
26	14	1799	G	OP1-P-OP2	5.03	127.15	119.60
26	1H	1923	U	C6-N1-C2	-5.03	117.98	121.00
27	16	100	G	N3-C4-N9	5.03	129.02	126.00
26	14	389	G	N3-C4-N9	5.03	129.02	126.00
26	14	463	G	O5'-P-OP1	5.03	116.74	110.70
26	14	1588	C	O5'-P-OP2	-5.03	101.17	105.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	606	U	N1-C2-N3	5.03	117.92	114.90
26	1H	1602	U	N1-C2-O2	-5.03	119.28	122.80
26	1H	1761	C	N3-C4-N4	5.03	121.52	118.00
26	1H	1898	U	OP1-P-O3'	5.03	116.27	105.20
26	1H	2382	G	C8-N9-C4	-5.03	104.39	106.40
34	58	134	ARG	N-CA-C	5.03	124.58	111.00
26	14	1776	G	C6-C5-N7	-5.03	127.38	130.40
1	13	335	C	C5-C6-N1	-5.03	118.49	121.00
1	13	578	C	N3-C4-N4	5.03	121.52	118.00
1	13	585	G	C8-N9-C4	5.03	108.41	106.40
1	13	1222	G	O5'-P-OP1	5.03	116.73	110.70
26	1H	470	A	C5-C6-N6	-5.03	119.68	123.70
26	1H	804	A	O4'-C1'-N9	5.03	112.22	108.20
26	14	322	A	N1-C6-N6	-5.03	115.58	118.60
26	14	1001	A	N9-C4-C5	5.03	107.81	105.80
26	14	1632	A	C5-C6-N6	-5.03	119.68	123.70
1	13	1054	C	O5'-P-OP2	5.03	116.73	110.70
26	1H	1463	C	C6-N1-C2	-5.03	118.29	120.30
26	1H	1777	U	OP2-P-O3'	5.03	116.26	105.20
26	1H	2035	G	O4'-C1'-N9	5.03	112.22	108.20
26	1H	2230	G	N3-C4-N9	-5.03	122.98	126.00
27	16	28	C	N3-C4-C5	-5.03	119.89	121.90
27	16	100	G	N3-C4-C5	-5.03	126.09	128.60
1	1G	360	A	C8-N9-C4	5.03	107.81	105.80
26	14	478	A	O5'-P-OP1	-5.03	101.17	105.70
26	14	2597	G	C5-C6-N1	-5.03	108.99	111.50
26	1H	673	C	C5-C4-N4	-5.03	116.68	120.20
26	1H	842	G	OP2-P-O3'	5.03	116.26	105.20
26	1H	2419	U	OP1-P-OP2	-5.03	112.06	119.60
31	41	34	LEU	CA-CB-CG	5.03	126.86	115.30
26	14	990	A	C8-N9-C4	-5.03	103.79	105.80
26	14	2059	A	C6-N1-C2	-5.03	115.58	118.60
26	14	2435	A	N7-C8-N9	5.03	116.31	113.80
26	1H	669	G	N7-C8-N9	5.02	115.61	113.10
26	1H	2672	G	C6-C5-N7	-5.02	127.39	130.40
1	1G	1480	G	C5-C6-O6	-5.02	125.58	128.60
26	14	1197	G	C5-C6-O6	5.02	131.61	128.60
26	14	1904	G	N1-C6-O6	-5.02	116.89	119.90
26	14	2615	U	N1-C2-O2	5.02	126.32	122.80
26	1H	212	G	OP2-P-O3'	5.02	116.25	105.20
26	1H	391	G	C8-N9-C1'	-5.02	120.47	127.00
26	1H	463	G	OP1-P-O3'	5.02	116.25	105.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	534	U	O5'-P-OP1	-5.02	101.18	105.70
26	1H	1368	G	C2-N3-C4	5.02	114.41	111.90
26	14	561	G	N3-C2-N2	-5.02	116.39	119.90
26	14	953	A	N1-C6-N6	5.02	121.61	118.60
26	14	2326	C	N3-C4-C5	-5.02	119.89	121.90
26	14	2330	G	C6-C5-N7	-5.02	127.39	130.40
1	13	557	G	N3-C2-N2	5.02	123.42	119.90
26	1H	842	G	C8-N9-C1'	5.02	133.53	127.00
26	1H	2530	A	C4-C5-N7	5.02	113.21	110.70
26	1H	2759	G	O4'-C1'-N9	-5.02	104.18	108.20
1	13	1299	A	C4-C5-N7	5.02	113.21	110.70
26	1H	456	C	C6-N1-C2	5.02	122.31	120.30
26	1H	1260	G	C4-C5-N7	-5.02	108.79	110.80
26	1H	1345	C	N3-C4-C5	5.02	123.91	121.90
26	1H	1368	G	OP1-P-OP2	5.02	127.13	119.60
26	1H	2446	G	C5-N7-C8	-5.02	101.79	104.30
26	1H	2589	A	N7-C8-N9	-5.02	111.29	113.80
26	14	2428	G	P-O3'-C3'	5.02	125.72	119.70
1	13	942	G	OP1-P-O3'	5.02	116.24	105.20
26	1H	196	A	C5-N7-C8	-5.02	101.39	103.90
26	1H	515	A	O4'-C1'-N9	5.02	112.22	108.20
26	1H	733	G	C5-C6-O6	5.02	131.61	128.60
26	1H	2401	U	C6-N1-C2	-5.02	117.99	121.00
28	11	43	ARG	NE-CZ-NH1	-5.02	117.79	120.30
26	14	787	U	OP1-P-OP2	-5.02	112.08	119.60
26	14	803	U	C5-C6-N1	-5.02	120.19	122.70
26	14	940	G	C2-N3-C4	5.02	114.41	111.90
26	14	1347	G	C5-C6-N1	-5.02	108.99	111.50
26	14	2442	C	OP2-P-O3'	5.02	116.24	105.20
26	1H	1334	G	N7-C8-N9	5.02	115.61	113.10
26	14	1950	G	N7-C8-N9	5.02	115.61	113.10
26	14	2335	A	O5'-P-OP1	-5.02	101.19	105.70
1	13	21	G	C4-C5-N7	-5.01	108.79	110.80
1	13	518	C	C2-N1-C1'	5.01	124.32	118.80
1	13	579	G	C5-C6-N1	-5.01	108.99	111.50
26	1H	120	U	C2-N3-C4	-5.01	123.99	127.00
26	1H	580	C	C4-C5-C6	5.01	119.91	117.40
26	1H	762	U	C6-N1-C1'	-5.01	114.18	121.20
26	1H	766	C	OP2-P-O3'	5.01	116.23	105.20
26	1H	2358	G	C5-C6-N1	5.01	114.01	111.50
33	61	110	ASP	C-N-CD	-5.01	109.57	120.60
26	14	330	A	C6-C5-N7	-5.01	128.79	132.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	766	C	C6-N1-C1'	5.01	126.82	120.80
26	14	1925	C	C2-N1-C1'	-5.01	113.28	118.80
26	14	2075	U	N3-C4-C5	5.01	117.61	114.60
26	14	2391	G	OP1-P-O3'	5.01	116.23	105.20
1	1G	1436	U	C5-C6-N1	5.01	125.21	122.70
1	13	1062	U	O5'-P-OP2	-5.01	101.19	105.70
1	13	1482	G	N1-C6-O6	5.01	122.91	119.90
26	1H	463	G	N1-C6-O6	-5.01	116.89	119.90
26	1H	1773	A	C8-N9-C4	5.01	107.80	105.80
1	1G	1432	G	C8-N9-C4	-5.01	104.39	106.40
26	14	819	A	N3-C4-C5	-5.01	123.29	126.80
26	14	1407	C	C6-N1-C2	-5.01	118.30	120.30
26	1H	138	G	N9-C1'-C2'	5.01	120.51	114.00
26	1H	232	G	N3-C4-N9	5.01	129.01	126.00
26	1H	829	A	OP1-P-OP2	5.01	127.11	119.60
26	1H	1060	U	P-O3'-C3'	5.01	125.71	119.70
26	1H	1783	A	C6-N1-C2	5.01	121.61	118.60
26	1H	2055	C	N1-C2-O2	-5.01	115.89	118.90
26	1H	2379	G	C8-N9-C4	5.01	108.40	106.40
26	1H	2774	C	C6-N1-C2	5.01	122.30	120.30
26	14	1543	A	O5'-P-OP1	5.01	116.71	110.70
26	14	2249	U	C2-N3-C4	5.01	130.00	127.00
26	1H	1951	U	N1-C2-O2	-5.01	119.30	122.80
1	1G	60	A	C8-N9-C4	5.01	107.80	105.80
1	1G	454	C	N3-C2-O2	-5.01	118.39	121.90
26	1H	457	A	C8-N9-C4	-5.01	103.80	105.80
26	1H	1799	G	N3-C4-C5	-5.01	126.10	128.60
26	1H	1816	G	C6-C5-N7	5.01	133.40	130.40
26	1H	2245	U	O4'-C1'-N1	5.01	112.20	108.20
26	1H	2443	C	C5-C4-N4	-5.01	116.69	120.20
23	2L	35	C	N3-C2-O2	-5.01	118.39	121.90
26	14	68	G	N1-C6-O6	5.01	122.90	119.90
26	14	465	G	C4-N9-C1'	5.01	133.01	126.50
26	14	759	G	OP1-P-OP2	-5.01	112.09	119.60
26	14	2831	G	C5-C6-O6	-5.01	125.60	128.60
26	1H	66	C	C5-C6-N1	5.00	123.50	121.00
27	16	81	G	C5-C6-O6	-5.00	125.60	128.60
26	14	1241	A	C5-C6-N1	-5.00	115.20	117.70
1	13	1221	G	OP2-P-O3'	5.00	116.21	105.20
26	1H	2712	U	P-O3'-C3'	5.00	125.70	119.70
26	14	265	A	C6-C5-N7	-5.00	128.80	132.30
26	14	1821	A	C5-C6-N6	-5.00	119.70	123.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	1983	C	C6-N1-C2	-5.00	118.30	120.30
26	14	2286	A	N7-C8-N9	5.00	116.30	113.80
26	14	2330	G	P-O3'-C3'	-5.00	113.70	119.70
1	13	22	G	OP2-P-O3'	5.00	116.20	105.20
1	13	1158	C	C6-N1-C1'	-5.00	114.80	120.80
26	1H	126	A	O5'-P-OP2	-5.00	101.20	105.70
26	1H	622	G	N3-C2-N2	5.00	123.40	119.90
26	1H	973	A	N1-C2-N3	5.00	131.80	129.30
26	1H	1256	G	C8-N9-C4	5.00	108.40	106.40
26	1H	1775	U	OP1-P-O3'	5.00	116.20	105.20
26	1H	2316	C	C6-N1-C2	-5.00	118.30	120.30
26	14	138	G	N1-C6-O6	5.00	122.90	119.90
26	14	574	C	C6-N1-C1'	5.00	126.80	120.80
26	14	651	G	C5-C6-O6	5.00	131.60	128.60
26	14	1371	G	N3-C2-N2	-5.00	116.40	119.90
26	14	1523	U	C5-C6-N1	5.00	125.20	122.70
26	14	2427	C	OP1-P-OP2	-5.00	112.10	119.60

There are no chirality outliers.

All (88) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
28	11	197	GLY	Peptide
28	11	237	GLU	Peptide
28	11	238	GLY	Peptide
28	11	47	GLY	Peptide
28	19	236	GLY	Peptide
28	19	237	GLU	Peptide
28	19	268	ARG	Peptide
28	19	32	SER	Peptide
28	19	37	LEU	Peptide
10	1A	55	LYS	Peptide
29	21	186	GLY	Peptide
29	21	56	PRO	Peptide
29	21	78	LEU	Peptide
35	25	47	ILE	Peptide
29	29	201	THR	Peptide
29	29	65	GLY	Peptide
29	29	72	VAL	Peptide
11	2I	100	ALA	Peptide
11	2I	101	SER	Peptide
36	35	110	TYR	Peptide

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Mol	Chain	Res	Type	Group
36	35	14	LYS	Peptide
36	35	36	LYS	Peptide
36	35	52	GLU	Peptide
36	35	64	LYS	Peptide
30	39	20	LEU	Peptide
30	39	85	GLY	Peptide
4	3E	166	LYS	Peptide
4	3E	30	LYS	Peptide
31	41	95	ARG	Peptide
37	45	134	ARG	Peptide
37	45	25	ASP	Peptide
37	45	86	GLY	Peptide
37	45	87	LYS	Peptide
13	4I	105	THR	Peptide
13	4I	107	ALA	Peptide
33	61	11	ASN	Peptide
33	61	115	ALA	Peptide
33	61	134	PRO	Peptide
33	61	82	ARG	Peptide
39	65	59	LYS	Peptide
33	69	112	LYS	Peptide
15	6I	22	THR	Peptide
40	75	104	ASN	Peptide
40	75	4	GLY	Peptide
16	7I	76	GLN	Peptide
41	85	98	LEU	Peptide
37	88	58	PHE	Peptide
37	88	78	PRO	Peptide
9	8E	110	GLU	Peptide
39	A8	107	GLU	Peptide
39	A8	89	ARG	Peptide
19	AI	6	LYS	Peptide
19	AI	7	LYS	Peptide
44	B5	24	GLY	Peptide
44	B5	61	GLY	Peptide
40	B8	58	ASN	Peptide
20	BA	101	GLY	Peptide
20	BA	11	SER	Peptide
45	C5	18	GLY	Peptide
45	C5	81	LYS	Peptide
45	C5	91	GLU	Peptide
41	C8	92	ARG	Peptide

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Mol	Chain	Res	Type	Group
46	D5	61	LEU	Peptide
43	E8	91	GLY	Peptide
48	F5	35	THR	Peptide
48	F5	85	LEU	Peptide
44	F8	67	GLY	Peptide
49	G5	15	LYS	Peptide
49	G5	17	SER	Peptide
49	G5	47	ASN	Peptide
45	G8	94	LYS	Peptide
46	H8	158	PRO	Peptide
47	I8	83	PRO	Peptide
47	I8	9	SER	Peptide
52	J5	3	LYS	Peptide
49	K8	17	SER	Peptide
49	K8	4	SER	Peptide
49	K8	46	GLN	Peptide
55	M5	40	GLU	Peptide
51	M8	38	LYS	Peptide
51	M8	45	GLY	Peptide
52	N8	58	LEU	Peptide
53	O8	25	LYS	Peptide
55	Q8	18	ALA	Peptide
55	Q8	19	SER	Peptide
55	Q8	27	THR	Peptide
55	Q8	29	LYS	Peptide
55	Q8	9	GLY	Peptide

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	13	32207	0	16254	745	0
1	1G	32182	0	16245	835	1
2	12	1924	0	1975	96	0
2	1E	1924	0	1975	96	0
3	22	1612	0	1677	71	0
3	2E	1605	0	1668	48	0
4	32	1702	0	1763	86	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	3E	1702	0	1762	81	0
5	42	1155	0	1213	60	0
5	4E	1155	0	1213	51	0
6	52	842	0	857	31	0
6	5E	842	0	857	21	0
7	62	1256	0	1296	53	0
7	6E	1256	0	1296	52	0
8	72	1115	0	1177	47	0
8	7E	1115	0	1177	59	0
9	82	1009	0	1037	68	0
9	8E	1009	0	1037	60	0
10	1A	801	0	849	44	0
10	1I	801	0	849	52	0
11	2A	864	0	881	32	0
11	2I	864	0	881	35	0
12	3A	975	0	1062	53	0
12	3I	975	0	1062	37	0
13	4A	933	0	992	64	0
13	4I	938	0	997	51	0
14	5A	475	0	511	30	0
14	5I	491	0	529	24	0
15	6A	733	0	771	29	0
15	6I	733	0	771	29	0
16	7A	705	0	725	25	0
16	7I	705	0	725	58	0
17	8A	834	0	904	22	0
17	8I	834	0	904	47	0
18	9A	590	0	662	32	0
18	9I	590	0	662	22	0
19	AA	624	0	636	40	0
19	AI	647	0	665	33	0
20	BA	762	0	861	34	0
20	BI	762	0	861	48	0
21	1B	217	0	234	15	0
21	1F	217	0	234	10	0
22	1K	1628	0	840	36	0
23	2K	1646	0	845	29	0
23	2L	1646	0	845	32	0
24	1L	1619	0	822	21	0
24	3K	1619	0	822	53	0
24	3L	1619	0	822	40	0
25	4K	281	0	142	6	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
25	4L	193	0	99	8	0
26	14	62647	0	31581	1303	0
26	1H	62707	0	31607	1456	1
27	16	2617	0	1328	55	0
27	1J	2617	0	1328	92	0
28	11	2115	0	2195	101	0
28	19	2120	0	2197	89	0
29	21	1568	0	1634	100	0
29	29	1568	0	1634	119	0
30	31	1585	0	1632	77	0
30	39	1627	0	1680	100	0
31	41	1473	0	1535	75	0
31	49	1473	0	1535	61	0
32	51	1336	0	1418	75	0
32	59	1299	0	1371	68	0
33	61	1136	0	1223	63	0
33	69	1136	0	1223	51	0
34	15	1104	0	1180	56	0
34	58	1104	0	1180	52	0
35	25	932	0	996	37	0
35	68	932	0	996	46	0
36	35	1144	0	1228	81	0
36	78	1144	0	1228	102	0
37	45	1121	0	1179	61	0
37	88	1086	0	1129	65	0
38	55	959	0	1021	47	0
38	98	967	0	1033	58	0
39	65	881	0	943	68	0
39	A8	881	0	943	61	0
40	75	1141	0	1202	69	0
40	B8	1141	0	1202	55	0
41	85	963	0	1022	49	0
41	C8	963	0	1022	62	0
42	95	778	0	852	62	0
42	D8	778	0	852	42	0
43	A5	899	0	964	31	0
43	E8	899	0	964	33	0
44	B5	725	0	778	34	0
44	F8	742	0	803	42	0
45	C5	794	0	884	56	0
45	G8	791	0	881	55	0
46	D5	1428	0	1454	69	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
46	H8	1397	0	1430	56	0
47	E5	606	0	628	39	0
47	I8	626	0	642	38	0
48	F5	762	0	848	30	0
48	J8	762	0	848	28	0
49	G5	558	0	610	27	0
49	K8	563	0	612	44	0
50	H5	468	0	518	17	0
50	L8	452	0	503	35	0
51	I5	515	0	514	33	0
51	M8	533	0	526	42	0
52	J5	434	0	454	23	0
52	N8	453	0	475	18	0
53	K5	389	0	404	18	0
53	O8	389	0	404	30	0
54	L5	398	0	441	19	0
54	P8	391	0	432	14	0
55	M5	477	0	540	44	0
55	Q8	485	0	551	72	0
56	11	2	0	0	0	0
56	13	152	0	0	0	0
56	14	426	0	0	0	0
56	15	1	0	0	0	0
56	16	15	0	0	0	0
56	1G	88	0	0	0	0
56	1H	506	0	0	0	0
56	1J	10	0	0	0	0
56	21	2	0	0	0	0
56	25	2	0	0	0	0
56	29	2	0	0	0	0
56	2K	7	0	0	0	0
56	2L	3	0	0	0	0
56	31	1	0	0	0	0
56	32	1	0	0	0	0
56	35	1	0	0	0	0
56	39	3	0	0	0	0
56	3I	2	0	0	0	0
56	41	2	0	0	0	0
56	45	2	0	0	0	0
56	49	1	0	0	0	0
56	4K	1	0	0	0	0
56	52	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
56	5E	1	0	0	0	0
56	6A	1	0	0	0	0
56	78	2	0	0	0	0
56	85	1	0	0	0	0
56	88	2	0	0	0	0
56	C5	1	0	0	0	0
56	E5	1	0	0	0	0
56	G8	1	0	0	0	0
56	I8	2	0	0	0	0
56	J8	1	0	0	0	0
56	L8	1	0	0	0	0
56	M5	1	0	0	0	0
56	P8	1	0	0	0	0
57	32	1	0	0	0	0
57	3E	1	0	0	0	0
57	5A	1	0	0	0	0
57	5I	1	0	0	0	0
57	C5	1	0	0	0	0
57	G8	1	0	0	0	0
58	11	9	0	0	3	0
58	13	164	0	0	12	0
58	14	543	0	0	112	0
58	15	1	0	0	0	0
58	16	6	0	0	2	0
58	19	3	0	0	3	0
58	1G	64	0	0	12	0
58	1H	920	0	0	187	0
58	1I	1	0	0	1	0
58	1J	18	0	0	2	0
58	21	5	0	0	2	0
58	29	3	0	0	0	0
58	31	7	0	0	0	0
58	32	1	0	0	0	0
58	35	1	0	0	0	0
58	39	8	0	0	0	0
58	3E	1	0	0	0	0
58	3I	1	0	0	0	0
58	4E	1	0	0	0	0
58	4K	1	0	0	0	0
58	5I	1	0	0	0	0
58	6I	1	0	0	0	0
58	78	6	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
58	7A	1	0	0	0	0
58	7I	1	0	0	0	0
58	85	1	0	0	0	0
58	B8	1	0	0	0	0
58	C8	2	0	0	0	0
58	E8	1	0	0	0	0
58	F8	1	0	0	0	0
58	G8	3	0	0	1	0
58	J8	1	0	0	0	0
58	L5	2	0	0	0	0
58	L8	3	0	0	1	0
58	P8	1	0	0	0	0
58	Q8	2	0	0	0	0
All	All	299951	0	200381	8425	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 17.

All (8425) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:1H:1604:C:OP2	58:1H:3643:HOH:O	1.58	1.18
40:B8:50:ILE:HD11	40:B8:102:ILE:HD11	1.35	1.08
26:1H:2576:G:OP1	58:1H:3756:HOH:O	1.73	1.07
26:1H:1614:A:OP1	58:1H:3859:HOH:O	1.76	1.03
26:14:1774:C:OP1	58:14:3564:HOH:O	1.77	1.03
26:1H:2308:G:H1	26:1H:2311:A:H2	1.04	1.03
26:1H:1783:A:OP2	58:1H:3629:HOH:O	1.75	1.02
26:1H:1771:C:HO2'	26:1H:1786:A:H8	1.03	1.02
26:14:2701:C:H3'	26:14:2702:U:H5''	1.42	1.01
26:1H:862:G:OP2	58:1H:3905:HOH:O	1.78	1.01
26:1H:574:C:OP1	58:1H:4144:HOH:O	1.77	1.01
1:1G:963:G:H21	10:1A:55:LYS:HE3	1.25	1.01
26:1H:1622:G:OP2	58:1H:4336:HOH:O	1.77	1.00
26:1H:1496:A:H8	26:1H:1577:C:HO2'	1.01	1.00
26:14:1828:G:OP2	58:14:3539:HOH:O	1.79	0.99
24:3L:71:G:O2'	26:14:1851:U:O2'	1.78	0.99
26:14:1582:C:HO2'	26:14:1586:A:H8	1.03	0.99
1:1G:588:G:H1	1:1G:651:C:H42	1.07	0.98
36:78:19:VAL:HG12	36:78:21:ARG:H	1.29	0.98
26:1H:155:C:H42	26:1H:171:G:H1	1.03	0.98

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:1H:741:G:OP1	58:1H:4205:HOH:O	1.80	0.97
22:1K:74:C:H42	26:1H:2554:U:H3	1.12	0.97
27:1J:15:A:H5'	27:1J:16:G:H8	1.30	0.97
26:1H:1899:G:H22	26:1H:1902:C:N4	1.63	0.97
26:14:676:A:H8	26:14:2069:G:H21	1.06	0.97
26:14:1496:A:H8	26:14:1577:C:HO2'	1.05	0.96
26:1H:2154:G:HO2'	26:1H:2155:G:H8	1.09	0.96
26:1H:567:A:OP1	58:1H:3604:HOH:O	1.80	0.96
26:1H:1780:A:OP1	58:1H:3625:HOH:O	1.81	0.96
26:1H:566:U:OP1	36:78:29:LYS:NZ	1.97	0.96
26:14:662:G:H5'	36:35:15:ARG:HA	1.48	0.95
26:14:1327:C:OP2	58:14:3606:HOH:O	1.82	0.95
24:1L:76:A:N7	26:14:2583:G:N2	2.13	0.95
26:1H:913:U:O4	58:1H:4214:HOH:O	1.84	0.95
26:1H:607:U:H3	26:1H:621:A:H2	1.14	0.95
26:1H:1828:G:OP2	58:1H:3705:HOH:O	1.83	0.95
26:1H:943:U:OP2	58:1H:4419:HOH:O	1.84	0.95
26:14:1359:A:H62	26:14:1372:U:H3	1.12	0.94
26:1H:511:U:OP2	58:1H:4189:HOH:O	1.84	0.94
2:1E:185:ILE:HG22	2:1E:199:TYR:HB2	1.46	0.94
26:14:1613:G:N7	58:14:3799:HOH:O	2.00	0.94
31:41:112:PRO:HB3	51:M8:37:SER:H	1.29	0.94
26:1H:330:A:H2	26:1H:1210:A:HO2'	1.13	0.94
26:1H:2017:U:OP1	58:1H:4151:HOH:O	1.84	0.93
26:14:2588:G:OP1	58:14:3588:HOH:O	1.87	0.93
1:13:200:G:H1	1:13:217:C:H42	1.14	0.93
1:1G:1277:C:HO2'	1:1G:1279:A:H8	1.05	0.93
26:1H:593:G:H4'	55:Q8:60:LEU:HD13	1.49	0.93
27:1J:80:U:H2'	27:1J:81:G:H21	1.33	0.93
26:1H:945:A:OP1	58:1H:3982:HOH:O	1.86	0.93
26:1H:2419:U:O4	55:Q8:29:LYS:NZ	2.02	0.92
24:3K:76:A:H8	26:1H:2394:C:H42	1.15	0.92
26:1H:1997:G:OP2	58:1H:3916:HOH:O	1.88	0.91
26:1H:1899:G:H22	26:1H:1902:C:H41	1.04	0.91
27:16:101:A:OP2	58:16:306:HOH:O	1.89	0.91
2:12:12:GLU:HB3	2:12:213:LEU:HD22	1.52	0.91
1:13:1368:G:H5''	9:8E:112:LYS:HB3	1.52	0.91
26:1H:2701:C:H3'	26:1H:2702:U:H5''	1.53	0.91
31:49:104:GLU:HG2	51:I5:23:GLU:HG2	1.51	0.91
1:13:664:G:H22	1:13:741:G:H1	1.18	0.91
26:14:1614:A:OP1	58:14:3513:HOH:O	1.88	0.91

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:741:G:OP1	58:14:3561:HOH:O	1.89	0.90
28:11:228:PRO:O	58:11:407:HOH:O	1.88	0.90
1:13:1502:A:H2	1:13:1505:G:H1	1.15	0.90
26:14:450:G:O6	58:14:3864:HOH:O	1.89	0.90
26:1H:1603:A:OP1	58:1H:3643:HOH:O	1.88	0.90
26:14:71:A:H2	44:B5:31:HIS:HE2	1.15	0.90
26:1H:249:C:OP1	58:1H:3665:HOH:O	1.89	0.90
26:1H:326:G:N7	58:1H:4427:HOH:O	2.05	0.90
26:1H:973:A:OP2	58:1H:3817:HOH:O	1.89	0.90
29:21:135:HIS:NE2	58:21:402:HOH:O	2.04	0.90
26:14:1416:G:H1	26:14:1582:C:H42	1.20	0.89
26:1H:568:U:O4	58:1H:3817:HOH:O	1.90	0.89
32:59:159:GLU:O	32:59:163:TYR:OH	1.88	0.89
1:1G:1502:A:H2	1:1G:1505:G:H1	1.21	0.89
45:G8:87:LYS:HB3	45:G8:94:LYS:HG2	1.55	0.89
26:14:323:G:HO2'	26:14:1205:U:H3	1.03	0.89
30:39:188:ARG:HA	36:35:3:LEU:HD11	1.52	0.89
1:1G:533:A:OP1	58:1G:1725:HOH:O	1.91	0.89
26:14:1899:G:H21	26:14:1902:C:N4	1.70	0.89
26:14:1757:U:H3	26:14:1762:A:H2	1.22	0.88
26:14:125:G:H5''	54:L5:19:ARG:HD3	1.55	0.88
39:A8:34:HIS:HB2	39:A8:36:TYR:HE1	1.39	0.88
26:14:567:A:OP1	58:14:3841:HOH:O	1.91	0.88
26:14:1043:C:N3	26:14:1112:G:N2	2.22	0.87
24:3K:72:C:H3'	24:3K:73:A:H5''	1.57	0.87
26:1H:1970:A:OP2	58:1H:4211:HOH:O	1.91	0.87
7:6E:111:ARG:NH1	7:6E:113:GLU:OE2	2.07	0.87
1:1G:589:C:N3	1:1G:650:G:N2	2.22	0.87
26:1H:1728:G:H8	26:1H:1732:A:H62	1.23	0.87
26:14:2357:U:OP1	47:E5:20:ARG:NH1	2.07	0.87
26:1H:674:G:OP2	58:1H:3846:HOH:O	1.93	0.87
1:1G:976:G:N2	1:1G:1362(A):C:OP2	2.07	0.86
26:14:1141:U:OP2	34:15:63:THR:OG1	1.92	0.86
26:1H:748:G:OP2	58:1H:4196:HOH:O	1.92	0.86
1:13:538:G:H5''	12:3I:114:LYS:HB2	1.55	0.86
43:A5:88:ARG:NH1	43:A5:94:ASP:OD2	2.09	0.86
2:12:91:PRO:HG3	2:12:154:LEU:HB2	1.55	0.86
26:1H:878:A:N6	26:1H:899:A:O2'	2.08	0.86
26:1H:510:C:OP1	58:1H:4189:HOH:O	1.93	0.86
26:1H:330:A:HO2'	26:1H:331:A:H8	1.21	0.86
1:1G:750:G:N2	15:6A:23:GLY:O	2.08	0.86

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:1372:U:OP2	58:14:3971:HOH:O	1.92	0.86
26:14:2035:G:OP1	58:14:3847:HOH:O	1.92	0.86
15:6I:17:ARG:HD3	15:6I:26:GLU:HG3	1.57	0.86
45:G8:84:ARG:NH2	58:G8:302:HOH:O	2.08	0.86
36:78:86:LYS:HB3	36:78:118:GLY:HA3	1.58	0.86
8:72:86:ILE:HG21	8:72:133:LEU:HD23	1.58	0.86
26:1H:392:C:OP1	58:1H:3721:HOH:O	1.94	0.86
1:1G:1154:G:H2'	1:1G:1155:G:H8	1.41	0.86
26:1H:2256:G:N7	58:1H:3987:HOH:O	2.09	0.86
36:35:19:VAL:HG13	36:35:21:ARG:H	1.40	0.85
26:1H:2033:A:OP1	58:1H:3948:HOH:O	1.92	0.85
2:1E:67:THR:HG21	2:1E:155:LEU:HG	1.59	0.85
26:1H:987:G:OP2	58:1H:3911:HOH:O	1.94	0.85
36:35:39:LYS:HD2	36:35:45:LEU:HD21	1.56	0.85
5:42:102:ALA:HB1	5:42:106:PRO:HG2	1.58	0.85
16:7A:40:ASP:HB3	16:7A:48:TRP:HB2	1.58	0.85
26:1H:945:A:OP1	58:1H:3979:HOH:O	1.93	0.85
1:13:452:A:OP1	16:7I:43:LYS:NZ	2.09	0.85
26:1H:1601:G:N7	58:1H:4227:HOH:O	2.09	0.85
26:1H:1479:G:N7	26:1H:1510:A:N6	2.24	0.85
26:1H:2053:G:OP2	58:1H:3755:HOH:O	1.92	0.85
1:13:1348:U:H3	1:13:1374:A:H2	1.23	0.85
11:2I:99:GLN:HG2	11:2I:105:VAL:HG11	1.59	0.85
40:75:64:ARG:HB2	40:75:73:GLU:HG2	1.59	0.85
26:1H:2593:U:O4	58:1H:3678:HOH:O	1.92	0.85
24:1L:74:C:N4	26:14:2554:U:O2	2.09	0.85
47:E5:53:MET:HG3	47:E5:59:LEU:HD23	1.58	0.85
26:1H:138:G:N2	44:F8:44:GLU:OE2	2.08	0.85
1:13:768:A:OP2	58:13:1890:HOH:O	1.94	0.84
26:1H:399:G:OP2	58:1H:4236:HOH:O	1.95	0.84
26:1H:49:A:N7	26:1H:120:U:H5	1.75	0.84
26:1H:1434:A:H61	26:1H:1558:A:N6	1.75	0.84
2:1E:120:ALA:O	2:1E:124:SER:OG	1.95	0.84
26:1H:1386:C:H2'	26:1H:1387:C:H6	1.39	0.84
26:14:1997:G:OP2	58:14:3782:HOH:O	1.94	0.84
42:95:71:LEU:H	42:95:86:GLY:HA2	1.41	0.84
26:1H:2582:G:OP2	58:1H:4163:HOH:O	1.94	0.84
26:1H:2597:G:O3'	58:1H:3656:HOH:O	1.96	0.84
26:14:2114:A:N6	26:14:2119:A:N7	2.26	0.84
26:14:138:G:N2	44:B5:44:GLU:OE2	2.09	0.84
36:78:47:ASP:OD2	36:78:50:ARG:NH2	2.11	0.84

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:58:47:ALA:HB2	34:58:112:LEU:HD11	1.57	0.84
55:Q8:49:VAL:HG22	55:Q8:50:LEU:H	1.43	0.84
26:1H:1525:G:H2'	26:1H:1526:G:H8	1.42	0.84
44:F8:67:GLY:O	44:F8:69:TYR:N	2.09	0.84
29:21:201:THR:HG22	29:21:203:LYS:H	1.41	0.84
26:1H:1581:G:N7	58:1H:4513:HOH:O	2.11	0.84
26:1H:1156:A:OP2	58:1H:3762:HOH:O	1.96	0.84
26:1H:2392:A:H2	26:1H:2424:C:H42	1.22	0.84
1:1G:680:C:H42	1:1G:710:G:H1	1.25	0.84
26:1H:751:A:OP1	58:1H:3861:HOH:O	1.96	0.83
24:3K:13:C:H2'	24:3K:14:A:H8	1.43	0.83
27:1J:15:A:H5'	27:1J:16:G:C8	2.13	0.83
1:1G:590:C:O2	1:1G:649:G:N2	2.11	0.83
46:H8:72:ARG:NH2	46:H8:97:GLU:O	2.10	0.83
26:1H:878:A:N6	26:1H:899:A:HO2'	1.75	0.83
26:1H:1899:G:N2	26:1H:1902:C:H41	1.76	0.83
42:95:35:LEU:HD12	42:95:37:VAL:HG11	1.61	0.83
42:95:69:LYS:HB3	42:95:86:GLY:HA3	1.59	0.83
26:1H:1265:A:OP2	58:1H:3612:HOH:O	1.95	0.83
26:14:1485:G:H1	26:14:1504:C:H42	1.27	0.83
26:1H:563:G:OP2	58:1H:3638:HOH:O	1.97	0.83
26:1H:2032:G:H21	29:21:146:THR:HG23	1.44	0.83
39:65:50:SER:O	39:65:76:LYS:NZ	2.10	0.83
1:13:1008:C:N4	1:13:1021:G:O6	2.11	0.83
26:14:2255:G:OP2	58:14:3652:HOH:O	1.95	0.83
26:1H:259:G:HO2'	26:1H:621:A:HO2'	1.23	0.83
1:1G:286:G:N7	58:1G:1753:HOH:O	2.11	0.83
26:1H:1665:A:OP2	58:1H:3712:HOH:O	1.95	0.83
26:1H:2502:G:OP2	58:1H:3635:HOH:O	1.95	0.83
26:1H:2127:G:H22	26:1H:2162:G:H1'	1.41	0.83
2:12:75:LYS:HA	2:12:78:GLN:HB2	1.59	0.82
1:1G:456:C:H42	1:1G:476:G:H1	1.27	0.82
26:1H:1516:U:H2'	26:1H:1517:G:H8	1.43	0.82
26:1H:71:A:H2	44:F8:31:HIS:HE2	1.26	0.82
26:14:1263:U:OP2	58:14:3977:HOH:O	1.96	0.82
29:29:101:ARG:HB2	29:29:203:LYS:HD3	1.60	0.82
26:14:517:C:OP1	52:J5:16:ARG:NH2	2.12	0.82
13:4A:66:LEU:HA	13:4A:70:LEU:HB2	1.61	0.82
26:1H:576:U:OP1	58:1H:3622:HOH:O	1.95	0.82
26:14:1970:A:OP2	58:14:3577:HOH:O	1.95	0.82
1:1G:998:G:N2	1:1G:1043:C:N3	2.27	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1G:54:C:N4	1:1G:353:A:OP2	2.12	0.82
26:1H:120:U:OP1	58:1H:4405:HOH:O	1.97	0.82
6:5E:50:TYR:OH	18:9I:74:ARG:O	1.96	0.82
26:14:84:A:N6	26:14:102:G:O2'	2.12	0.82
20:BI:71:THR:HG22	20:BI:72:LEU:H	1.45	0.82
1:13:156:G:H1	1:13:165:C:H42	1.24	0.82
48:J8:41:ARG:HG3	48:J8:41:ARG:HH11	1.41	0.82
29:21:57:LYS:HG3	29:21:59:VAL:HG12	1.61	0.82
44:B5:8:ILE:O	49:G5:36:ARG:NH2	2.12	0.82
50:L8:37:LEU:HD12	50:L8:43:ILE:HD13	1.61	0.82
26:1H:2062:A:OP1	58:1H:3788:HOH:O	1.97	0.82
26:14:2074:U:OP1	58:14:3511:HOH:O	1.96	0.82
26:14:586:A:OP2	58:14:3535:HOH:O	1.96	0.81
22:1K:22:G:O6	22:1K:46:7MG:N2	2.12	0.81
55:Q8:14:VAL:HG21	55:Q8:21:LYS:HZ2	1.43	0.81
26:1H:676:A:H8	26:1H:2069:G:H21	1.28	0.81
32:51:124:GLU:HB3	32:51:132:ARG:HB3	1.62	0.81
24:1L:50:U:O4	24:1L:64:A:N6	2.13	0.81
37:88:109:VAL:HG13	37:88:113:GLN:HB3	1.61	0.81
1:13:737:A:H2'	1:13:738:C:H6	1.44	0.81
45:G8:100:ALA:HB1	45:G8:101:LYS:HB2	1.62	0.81
27:1J:18:G:H1	27:1J:65:C:H42	1.23	0.81
26:14:2343:C:O2'	26:14:2373:G:O2'	1.96	0.81
26:1H:2608:G:N7	58:1H:4164:HOH:O	2.13	0.81
1:13:1007:C:H42	1:13:1022:G:H1	1.28	0.81
35:25:115:VAL:HG13	35:25:121:VAL:HG21	1.62	0.81
36:78:138:LEU:HD12	36:78:144:GLU:HG3	1.61	0.81
26:14:1622:G:OP2	58:14:3933:HOH:O	1.96	0.81
40:B8:21:GLU:OE1	40:B8:91:ARG:NH2	2.13	0.81
26:14:2415:G:H4'	36:35:67:MET:H	1.44	0.81
14:5A:21:TYR:OH	14:5A:23:ARG:NH2	2.14	0.81
26:1H:620:G:H4'	26:1H:621:A:H5''	1.63	0.81
30:39:25:PRO:HD3	30:39:115:ALA:HB1	1.61	0.81
3:22:182:ILE:HG22	3:22:203:PHE:HA	1.63	0.81
1:13:559:A:OP1	5:4E:126:ARG:NH2	2.13	0.81
26:14:1754:C:OP1	40:75:96:ARG:NH1	2.14	0.81
26:1H:142:G:H1'	44:F8:37:THR:HG21	1.60	0.81
1:1G:21:G:OP1	58:1G:1737:HOH:O	1.99	0.81
26:1H:453:C:OP1	58:1H:3839:HOH:O	1.99	0.81
26:1H:1395:A:OP2	58:1H:3645:HOH:O	1.97	0.81
26:1H:1678:G:N2	26:1H:1989:G:H22	1.78	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:6E:111:ARG:HD2	7:6E:123:GLU:HB2	1.63	0.80
26:1H:442:G:H1'	30:31:48:THR:HG21	1.63	0.80
26:1H:860:U:H5	26:1H:917:A:C2	1.99	0.80
32:59:10:PRO:HD2	32:59:50:VAL:HG13	1.62	0.80
28:11:85:ASP:HB2	28:11:92:ILE:HG12	1.62	0.80
33:61:7:GLU:HA	33:61:15:VAL:HG22	1.63	0.80
26:1H:2035:G:OP1	58:1H:3744:HOH:O	1.99	0.80
26:1H:654(E):C:N3	26:1H:654(P):G:N2	2.29	0.80
45:C5:17:SER:OG	45:C5:18:GLY:N	2.13	0.80
1:1G:1298:C:OP2	7:62:114:ARG:NH2	2.14	0.80
26:1H:1187:G:OP2	58:1H:3824:HOH:O	1.98	0.80
26:1H:780:G:H21	26:1H:783:A:H62	1.28	0.80
53:K5:33:LYS:HE2	53:K5:34:LEU:H	1.46	0.80
26:1H:654(G):C:O2	26:1H:654(N):G:N2	2.15	0.80
34:15:67:LEU:HG	34:15:88:GLU:HG2	1.62	0.80
11:2A:54:ARG:NH2	24:3L:39:U:O2'	2.12	0.80
1:13:310:G:OP2	16:7I:27:LYS:NZ	2.13	0.80
26:14:574:C:OP2	58:14:3825:HOH:O	1.99	0.80
26:14:2304:G:N2	26:14:2312:U:O4	2.14	0.80
26:1H:192:C:OP2	58:1H:4128:HOH:O	2.00	0.80
4:32:30:LYS:HB3	4:32:35:ARG:HD2	1.64	0.80
8:7E:10:LEU:HD22	8:7E:83:ILE:HD11	1.63	0.80
26:1H:1399:C:N4	58:1H:4305:HOH:O	2.14	0.80
26:1H:2577:A:OP1	58:1H:3752:HOH:O	1.99	0.80
20:BI:73:HIS:HB3	20:BI:74:LYS:HD2	1.64	0.80
1:1G:617:G:H1	1:1G:623:C:H42	1.28	0.80
26:14:1249:U:OP1	58:14:3530:HOH:O	2.00	0.79
26:14:800:A:OP1	58:14:3859:HOH:O	1.98	0.79
26:14:739:G:OP1	58:14:3506:HOH:O	1.99	0.79
4:3E:15:GLU:OE1	4:3E:66:ARG:NH1	2.16	0.79
33:61:110:ASP:HB2	33:61:112:LYS:HG2	1.62	0.79
5:4E:8:GLU:HG2	5:4E:34:VAL:HG22	1.65	0.79
47:E5:49:LYS:HG3	47:E5:80:HIS:HB3	1.63	0.79
26:1H:120:U:OP2	58:1H:3999:HOH:O	2.00	0.79
9:8E:21:PRO:HA	9:8E:59:PHE:HA	1.65	0.79
1:1G:961:U:O2	1:1G:1201:A:N6	2.15	0.79
1:13:1160:G:H1	1:13:1177:G:H22	1.30	0.79
27:16:12:C:O2	47:I8:74:ARG:NH1	2.15	0.79
26:1H:1316:U:H2'	26:1H:1317:A:H8	1.45	0.79
1:1G:1435:G:H2'	1:1G:1436:U:C6	2.17	0.79
26:14:2681:C:H5	26:14:2725:A:H62	1.26	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:848:G:H2'	26:14:849:A:C8	2.17	0.79
10:1I:77:PRO:HB2	10:1I:79:ARG:HH12	1.46	0.79
39:65:36:TYR:HA	39:65:52:SER:HB3	1.63	0.79
2:1E:189:ASP:HB2	2:1E:205:ASP:HB3	1.65	0.79
4:32:31:CYS:HB3	4:32:33:MET:HB2	1.65	0.79
26:14:751:A:OP1	58:14:3512:HOH:O	2.01	0.79
2:12:73:THR:HG21	2:12:97:TRP:H	1.48	0.79
1:1G:1069:C:O2'	1:1G:1192:C:O2	2.01	0.79
30:39:46:ARG:HG2	30:39:46:ARG:HH11	1.48	0.79
26:1H:879:G:O6	26:1H:898:C:N4	2.15	0.79
26:14:631:A:OP2	55:M5:47:LYS:NZ	2.11	0.79
2:1E:21:ARG:HB2	2:1E:39:ILE:HA	1.65	0.79
55:M5:40:GLU:HA	55:M5:43:GLN:HB2	1.65	0.79
30:39:117:ARG:HH12	36:35:1:MET:H2	1.30	0.79
27:1J:3:C:N3	27:1J:117:G:N2	2.30	0.79
45:C5:39:VAL:HG23	45:C5:41:GLY:H	1.48	0.79
26:14:2012:G:OP1	43:A5:11:ARG:NH2	2.16	0.79
27:1J:38:C:N3	27:1J:44:G:N2	2.29	0.79
26:1H:2046:G:H5'	52:N8:19:ARG:HA	1.63	0.78
26:1H:2056:G:OP2	58:1H:3617:HOH:O	2.01	0.78
26:14:833:U:O2	36:35:55:ARG:NH1	2.17	0.78
26:1H:654(H):G:N7	26:1H:654(N):G:N2	2.30	0.78
27:1J:38:C:H42	27:1J:44:G:H1	1.29	0.78
4:32:153:ARG:NH1	4:32:181:MET:SD	2.56	0.78
1:13:1213:A:O2'	1:13:1215:G:N7	2.14	0.78
26:1H:2126:A:N6	26:1H:2163:C:O2	2.16	0.78
27:16:100:G:OP1	58:16:303:HOH:O	2.01	0.78
10:1I:61:GLU:OE2	14:5I:45:ARG:NH1	2.16	0.78
39:A8:28:VAL:HG11	39:A8:98:VAL:HG13	1.65	0.78
1:1G:362:G:H4'	12:3A:33:ARG:HH21	1.49	0.78
26:1H:1633:G:OP2	58:1H:4369:HOH:O	2.01	0.78
1:13:972:C:OP1	58:13:1882:HOH:O	1.99	0.78
1:1G:376:G:H5''	16:7A:5:ARG:HD2	1.63	0.78
2:1E:100:GLY:O	2:1E:104:ASN:N	2.16	0.78
26:1H:2135:A:N6	26:1H:2156:G:O2'	2.17	0.78
26:1H:1778:U:H2'	26:1H:1784:A:N6	1.98	0.78
26:1H:1639:U:OP1	58:1H:3674:HOH:O	2.02	0.78
26:1H:2210:G:H3'	26:1H:2211:G:C8	2.19	0.78
9:8E:121:ARG:NH1	9:8E:122:ALA:O	2.16	0.78
26:14:2270:G:OP2	58:14:3728:HOH:O	2.01	0.78
26:14:1019:U:H3	26:14:1142(A):A:H62	1.30	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:1H:2334:G:O6	47:I8:74:ARG:NH2	2.15	0.78
31:41:66:GLN:OE1	31:41:98:ARG:NH1	2.15	0.78
36:78:2:LYS:HE3	36:78:4:SER:HB2	1.65	0.78
10:1I:57:LYS:HD2	10:1I:60:ARG:HH12	1.49	0.78
51:M8:48:ARG:HH11	51:M8:48:ARG:HA	1.49	0.78
29:29:66:HIS:NE2	29:29:73:GLU:OE1	2.15	0.78
26:14:2134:A:O2'	26:14:2159:G:N2	2.16	0.78
39:65:74:ALA:HB1	39:65:107:GLU:HB2	1.65	0.78
1:1G:490:G:OP2	4:32:132:ARG:NH2	2.16	0.78
46:H8:19:ARG:NH1	46:H8:84:GLU:O	2.17	0.78
26:14:2392:A:H2	26:14:2424:C:H42	1.29	0.78
28:11:93:ALA:HB3	28:11:105:ILE:HG22	1.64	0.78
26:14:593:G:H4'	55:M5:60:LEU:HD22	1.64	0.78
26:14:780:G:H21	26:14:783:A:H62	1.31	0.78
36:35:47:ASP:OD1	36:35:50:ARG:NH1	2.17	0.78
26:1H:1900:A:H5'	26:1H:1900:A:H8	1.47	0.78
12:3A:52:LEU:O	12:3A:54:LYS:NZ	2.17	0.78
19:AI:40:ILE:HG23	19:AI:41:VAL:HG13	1.64	0.78
1:1G:1002:G:N2	1:1G:1038:C:N3	2.32	0.78
1:1G:853:G:H2'	1:1G:854:G:H8	1.47	0.78
32:59:6:ARG:H	32:59:6:ARG:HH11	1.32	0.78
26:1H:1774:C:OP1	58:1H:3750:HOH:O	2.01	0.78
36:35:81:GLN:NE2	36:35:106:LEU:O	2.17	0.78
19:AA:41:VAL:HG12	19:AA:43:GLU:H	1.49	0.78
41:C8:8:VAL:HG23	41:C8:11:ARG:HH21	1.49	0.78
26:14:1702:G:N7	58:14:4006:HOH:O	2.16	0.78
26:1H:2789:C:O2	26:1H:2894:G:N2	2.15	0.78
33:69:130:TYR:HB3	33:69:136:VAL:HG13	1.66	0.78
26:1H:1287:A:N7	38:98:107:ASP:HB2	1.98	0.78
41:85:28:ARG:NH1	41:85:38:THR:OG1	2.17	0.77
26:14:1786:A:OP1	58:14:3569:HOH:O	2.02	0.77
1:1G:673:G:H2'	1:1G:674:G:C8	2.19	0.77
26:14:273(C):C:H42	26:14:363(C):G:H1	1.32	0.77
1:13:510:A:OP2	4:3E:49:ARG:NH2	2.17	0.77
26:14:270(Q):C:H5''	33:69:45:LYS:HE3	1.65	0.77
26:14:2352:A:C2	47:E5:33:ALA:HB1	2.17	0.77
26:14:1113:U:H5'	32:59:2:SER:HA	1.65	0.77
1:13:376:G:H1	1:13:387:U:H3	1.30	0.77
9:82:13:ALA:HB2	9:82:68:GLY:HA3	1.67	0.77
29:29:81:ILE:HG22	29:29:82:ARG:H	1.49	0.77
48:J8:65:SER:OG	48:J8:66:HIS:ND1	2.17	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1G:1210:C:HO2'	1:1G:1213:A:HO2'	1.29	0.77
46:D5:157:LEU:HB3	46:D5:161:VAL:HG11	1.65	0.77
52:N8:33:CYS:HB2	52:N8:40:LYS:HD3	1.65	0.77
35:25:13:ASN:HD21	35:25:96:THR:HG23	1.50	0.77
26:1H:509:C:O3'	58:1H:4191:HOH:O	2.02	0.77
26:14:654(D):G:H22	26:14:654(Q):C:H42	1.33	0.77
1:13:1029:G:H1'	1:13:1032(A):G:H22	1.49	0.77
26:1H:259:G:H21	26:1H:621:A:H8	1.30	0.77
45:C5:17:SER:HB2	45:C5:71:LYS:HD2	1.66	0.77
14:5A:45:ARG:O	14:5A:49:HIS:ND1	2.18	0.77
26:14:397:G:N7	58:14:3752:HOH:O	2.18	0.77
1:13:859:A:H2'	1:13:860:A:H8	1.49	0.77
2:12:47:THR:HG23	2:12:202:PRO:HG2	1.64	0.77
34:15:18:ALA:HA	34:15:21:LYS:HG3	1.66	0.77
33:61:144:VAL:HG13	33:61:145:VAL:HG23	1.65	0.77
27:1J:104:A:OP1	46:D5:72:ARG:NH1	2.18	0.77
1:1G:156:G:N2	1:1G:165:C:O2	2.18	0.77
42:95:71:LEU:N	42:95:86:GLY:HA2	1.99	0.77
26:1H:1782:C:OP1	58:1H:3627:HOH:O	2.02	0.77
26:14:920:G:H2'	26:14:921:G:H8	1.49	0.77
26:14:2068:U:H3	26:14:2430:A:H2	1.33	0.76
33:61:39:ALA:HB1	33:61:44:LEU:HD13	1.66	0.76
9:8E:50:LEU:HD23	9:8E:85:LEU:HD11	1.67	0.76
41:C8:92:ARG:O	41:C8:94:ASN:N	2.19	0.76
53:O8:41:PRO:HB2	53:O8:43:CYS:HB2	1.65	0.76
26:1H:751:A:OP1	58:1H:3860:HOH:O	2.03	0.76
26:14:2689:U:P	26:14:2719:G:H22	2.09	0.76
26:1H:187:G:N7	58:1H:4055:HOH:O	2.17	0.76
26:14:1992:G:N7	58:14:3597:HOH:O	2.17	0.76
1:13:812:C:N3	58:13:1802:HOH:O	2.18	0.76
47:E5:27:GLU:HG3	47:E5:68:GLU:HA	1.68	0.76
8:7E:7:ALA:HB2	8:7E:85:ARG:HH11	1.50	0.76
46:D5:4:ARG:HA	46:D5:58:VAL:HB	1.65	0.76
49:K8:15:LYS:H	49:K8:67:LYS:NZ	1.84	0.76
26:1H:192:C:N3	58:1H:3654:HOH:O	2.19	0.76
32:51:83:TYR:HD1	32:51:84:SER:H	1.31	0.76
48:J8:91:LYS:O	48:J8:94:LEU:N	2.18	0.76
1:1G:1359:C:O2'	1:1G:1361:G:N7	2.17	0.76
1:13:737:A:H2'	1:13:738:C:C6	2.20	0.76
1:13:1305:G:N2	1:13:1331:G:H2'	2.01	0.76
26:1H:2807:G:N1	26:1H:2893:G:O6	2.19	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:2419:U:O4	55:M5:31:HIS:ND1	2.19	0.76
1:1G:11:G:N2	1:1G:525:C:O2'	2.19	0.76
26:1H:1165:U:H2'	26:1H:1166:C:C6	2.20	0.76
26:1H:1814:G:O6	58:1H:4231:HOH:O	2.03	0.76
26:1H:1828:G:OP1	58:1H:3896:HOH:O	2.04	0.76
26:1H:330:A:O2'	26:1H:331:A:H8	1.68	0.76
26:14:2686:G:N7	58:14:3783:HOH:O	2.19	0.76
19:AA:50:ALA:HB1	19:AA:57:HIS:HB3	1.67	0.76
49:G5:25:VAL:HG12	49:G5:60:LEU:HD23	1.67	0.76
39:A8:93:LYS:HG2	39:A8:95:HIS:HB2	1.67	0.76
46:H8:76:LEU:HD22	46:H8:76:LEU:H	1.51	0.75
26:1H:450:G:OP2	58:1H:3835:HOH:O	2.04	0.75
34:15:47:ALA:HB2	34:15:112:LEU:HD21	1.66	0.75
32:59:125:VAL:HG22	32:59:126:PRO:HA	1.67	0.75
1:13:187:C:O2	1:13:191(A):G:N1	2.19	0.75
32:59:15:VAL:HG11	32:59:29:PRO:HD3	1.67	0.75
26:14:1970:A:OP1	58:14:3572:HOH:O	2.04	0.75
26:14:67:U:H3	26:14:74:A:H2	1.31	0.75
39:A8:37:ALA:HB2	39:A8:101:LEU:HD21	1.66	0.75
36:35:128:HIS:HA	36:35:147:LEU:HA	1.69	0.75
26:1H:2838:G:N7	58:1H:4314:HOH:O	2.19	0.75
26:14:958:U:OP2	37:45:14:ARG:NH1	2.15	0.75
42:95:80:GLN:HG3	42:95:81:TYR:H	1.51	0.75
1:13:1003:G:N2	1:13:1037:C:O2	2.19	0.75
26:14:1434:A:H61	26:14:1558:A:H62	1.33	0.75
1:1G:192:U:H2'	1:1G:193:C:H6	1.51	0.75
44:B5:41:ASN:HA	44:B5:44:GLU:HB2	1.69	0.75
27:1J:18:G:N2	27:1J:65:C:N3	2.35	0.75
26:14:1864:U:OP1	26:14:2410:G:O2'	2.03	0.75
33:61:37:VAL:HG12	33:61:38:LEU:HD12	1.68	0.75
26:14:2475:C:HO2'	26:14:2477:C:H5	1.35	0.75
1:1G:41:G:H2'	1:1G:42:G:C8	2.21	0.75
26:1H:2656:U:H3	26:1H:2665:A:H2	1.32	0.75
26:1H:2199:A:H5''	26:1H:2205:C:H5	1.52	0.75
1:1G:377:G:H1	1:1G:386:C:H42	1.33	0.75
15:6A:54:ARG:NH1	15:6A:58:MET:SD	2.59	0.75
26:1H:1495:A:OP2	58:1H:4102:HOH:O	2.04	0.75
10:1I:57:LYS:O	10:1I:60:ARG:NH1	2.19	0.75
26:14:2210:G:H3'	26:14:2211:G:C2	2.22	0.75
45:C5:97:ARG:NH2	45:C5:104:GLY:O	2.20	0.75
26:1H:2577:A:OP1	58:1H:3756:HOH:O	2.04	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:15:94:HIS:HB2	34:15:97:ARG:HG3	1.69	0.75
26:1H:1664:A:OP1	58:1H:3712:HOH:O	2.05	0.74
26:14:749:C:OP2	58:14:3643:HOH:O	2.03	0.74
24:3L:26:A:H61	24:3L:44:G:H1	1.31	0.74
29:21:105:THR:OG1	29:21:199:ARG:NH2	2.20	0.74
16:7I:4:ILE:HB	16:7I:66:PRO:HB3	1.69	0.74
26:1H:1580:A:OP1	58:1H:4512:HOH:O	2.04	0.74
27:1J:76:G:N7	58:1J:310:HOH:O	2.20	0.74
1:13:1086:U:O4	1:13:1099:G:N2	2.17	0.74
1:13:1004:A:O5'	1:13:1025:U:N3	2.15	0.74
26:14:2250:G:C4	37:45:82:ARG:HG3	2.23	0.74
40:B8:6:LEU:HA	40:B8:9:LEU:HB2	1.68	0.74
43:A5:65:LEU:HD13	43:A5:68:ARG:HD3	1.69	0.74
1:13:352:C:O2'	1:13:354:G:OP1	2.04	0.74
10:1I:34:VAL:HG12	10:1I:74:ILE:HG23	1.70	0.74
1:1G:587:G:N2	1:1G:754:C:OP2	2.20	0.74
1:1G:1248:A:N6	1:1G:1288:A:OP2	2.21	0.74
1:13:827:U:H5	1:13:872:A:N1	1.85	0.74
12:3I:52:LEU:O	12:3I:54:LYS:NZ	2.20	0.74
2:1E:33:TYR:HB2	2:1E:43:ASP:HB2	1.68	0.74
1:1G:1248:A:N3	9:82:70:LYS:NZ	2.33	0.74
29:29:16:ARG:NH2	29:29:171:GLU:OE1	2.20	0.74
26:14:2839:G:H5'	38:55:46:GLY:HA2	1.68	0.74
40:75:77:PRO:HG2	40:75:80:SER:HB2	1.69	0.74
26:1H:1791:A:H5'	28:11:206:LEU:HD12	1.68	0.74
26:1H:1611:C:OP2	58:1H:4174:HOH:O	2.05	0.74
24:3K:15:G:H1	24:3K:48:C:H41	1.34	0.74
26:1H:1900:A:C8	26:1H:1900:A:H5'	2.23	0.74
12:3I:126:LYS:HG3	12:3I:128:ALA:H	1.53	0.74
1:13:353:A:H5'	1:13:353:A:H8	1.52	0.74
1:13:766:A:OP2	58:13:1803:HOH:O	2.04	0.74
1:13:1149:C:H2'	1:13:1150:U:H6	1.52	0.74
26:14:2331:G:O3'	47:E5:43:THR:HG22	1.87	0.74
42:D8:60:GLU:HB2	42:D8:97:LYS:HE2	1.69	0.74
28:19:69:ARG:HD3	28:19:105:ILE:HD11	1.70	0.74
1:1G:1003:G:N2	1:1G:1037:C:O2	2.19	0.74
26:14:2873:A:H8	38:55:6:SER:H	1.35	0.74
47:I8:49:LYS:NZ	47:I8:68:GLU:OE2	2.13	0.74
26:1H:2059:A:OP2	58:1H:4045:HOH:O	2.04	0.74
4:32:98:GLU:OE2	4:32:103:ASN:ND2	2.18	0.74
26:14:1061:U:H4'	26:14:1070:A:H1'	1.68	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:1H:577:G:O6	58:1H:3922:HOH:O	2.06	0.74
1:1G:1002:G:H2'	1:1G:1003:G:C8	2.23	0.74
1:1G:1004:A:OP1	1:1G:1024:G:N1	2.17	0.74
26:1H:2583:G:OP2	58:1H:4163:HOH:O	2.05	0.74
13:4I:3:ARG:HE	13:4I:9:ILE:HD11	1.53	0.74
26:1H:1533:C:H3'	26:1H:1534:G:H5''	1.69	0.74
2:12:185:ILE:HG22	2:12:199:TYR:HB2	1.69	0.74
2:12:134:GLU:OE1	2:12:137:ARG:NH1	2.21	0.74
26:14:760:G:OP1	58:14:3870:HOH:O	2.05	0.74
1:13:1423:G:OP1	35:68:49:ARG:NH2	2.21	0.74
2:12:67:THR:HG21	2:12:155:LEU:HG	1.68	0.73
26:14:920:G:H2'	26:14:921:G:C8	2.23	0.73
1:13:1329:A:H5'	13:4I:29:ARG:HD2	1.68	0.73
26:14:877:U:O4	26:14:899:A:N6	2.21	0.73
1:1G:591:U:OP2	8:72:30:ARG:NH1	2.19	0.73
26:14:1434:A:H61	26:14:1558:A:N6	1.86	0.73
26:1H:1466:G:N2	26:1H:1547:C:N3	2.36	0.73
26:14:1041:C:H42	26:14:1114:G:H1	1.32	0.73
30:31:11:VAL:HG22	30:31:125:LEU:HB2	1.71	0.73
26:14:1856:G:H1	26:14:1886:C:H42	1.33	0.73
29:21:36:ARG:NH2	29:21:88:GLY:O	2.21	0.73
1:1G:1126:U:H4'	1:1G:1127:G:C8	2.23	0.73
26:14:751:A:OP1	58:14:3515:HOH:O	2.06	0.73
1:1G:353:A:H8	1:1G:353:A:H5'	1.53	0.73
1:13:963:G:H21	10:1I:55:LYS:HE2	1.53	0.73
26:1H:2199:A:H5'	26:1H:2205:C:OP2	1.89	0.73
37:88:51:ARG:HH12	37:88:52:VAL:HG23	1.52	0.73
1:13:74:C:H42	1:13:96:G:H1	1.36	0.73
26:1H:1249:U:OP1	58:1H:3828:HOH:O	2.06	0.73
26:1H:1434:A:H61	26:1H:1558:A:H61	1.34	0.73
26:14:1382:G:N7	58:14:3672:HOH:O	2.20	0.73
38:55:33:ARG:HD3	38:55:113:LEU:HD11	1.70	0.73
29:29:151:TYR:HD2	29:29:154:LYS:HZ2	1.35	0.73
32:51:4:ILE:HB	32:51:6:ARG:HG3	1.71	0.73
26:14:1900:A:OP2	58:14:3578:HOH:O	2.05	0.73
55:M5:52:LYS:HZ1	55:M5:53:PRO:HA	1.53	0.73
26:14:1270:C:H5''	26:14:1271:G:H5'	1.71	0.73
26:14:453:C:OP1	58:14:3860:HOH:O	2.06	0.73
36:78:50:ARG:HG3	55:Q8:58:ILE:HD11	1.69	0.73
49:K8:17:SER:HB3	49:K8:67:LYS:HE3	1.71	0.73
29:21:119:ARG:HD3	29:21:160:TYR:HB2	1.70	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
27:16:40:U:H1'	27:16:45:A:H61	1.52	0.73
1:13:177:C:OP1	20:BI:65:LYS:NZ	2.17	0.73
5:42:101:ILE:HD11	5:42:119:LEU:HD23	1.71	0.73
43:A5:13:SER:HB3	43:A5:16:LYS:HD2	1.70	0.73
16:7A:53:VAL:HG13	16:7A:79:VAL:HG22	1.71	0.73
26:1H:500:G:N7	58:1H:4351:HOH:O	2.22	0.73
30:31:103:LYS:HA	30:31:106:ARG:HG3	1.70	0.73
26:14:1249:U:OP1	58:14:3534:HOH:O	2.06	0.73
26:14:607:U:H3	26:14:621:A:H2	1.36	0.73
37:88:116:GLU:OE2	37:88:119:ARG:NH2	2.21	0.73
41:85:92:ARG:HH22	42:95:10:LYS:HA	1.53	0.73
26:1H:1525:G:H2'	26:1H:1526:G:C8	2.23	0.73
26:1H:1689:A:H62	26:1H:1698:A:H2	1.37	0.73
34:58:56:ASN:N	34:58:125:GLY:O	2.13	0.73
1:13:606:G:N3	1:13:632:A:N6	2.36	0.73
30:39:122:LYS:O	30:39:124:LEU:N	2.22	0.73
1:13:975:A:H4'	1:13:976:G:H5''	1.71	0.73
11:2A:85:ARG:HD3	11:2A:113:PRO:HD3	1.70	0.73
55:Q8:53:PRO:HA	55:Q8:55:ALA:N	2.04	0.72
19:AI:41:VAL:HG21	19:AI:67:VAL:HG12	1.68	0.72
26:1H:1441:G:H2'	26:1H:1442:G:H8	1.53	0.72
3:22:88:ARG:HB2	3:22:101:LEU:HD22	1.71	0.72
3:22:91:LEU:HB2	3:22:99:VAL:HG11	1.69	0.72
26:14:399:G:OP2	58:14:3682:HOH:O	2.06	0.72
26:1H:598:G:H5'	36:78:11:GLY:HA3	1.71	0.72
5:4E:137:GLU:OE1	5:4E:141:GLN:NE2	2.21	0.72
26:1H:574:C:OP2	58:1H:3919:HOH:O	2.07	0.72
26:1H:2062:A:N3	26:1H:2062:A:H2'	2.03	0.72
26:14:574:C:OP2	58:14:3827:HOH:O	2.07	0.72
55:M5:52:LYS:NZ	55:M5:53:PRO:HA	2.04	0.72
27:16:42:C:O3'	31:41:67:LYS:NZ	2.22	0.72
16:7I:53:VAL:HG13	16:7I:79:VAL:HG22	1.70	0.72
37:88:35:VAL:HG13	37:88:130:LYS:HB3	1.72	0.72
26:1H:1797:C:HO2'	28:11:259:THR:HG1	1.35	0.72
13:4A:97:PRO:HA	13:4A:110:ARG:HD3	1.72	0.72
24:3K:30:G:H1	24:3K:40:C:H42	1.37	0.72
26:1H:1828:G:OP1	58:1H:3893:HOH:O	2.06	0.72
1:1G:888:G:O2'	1:1G:1488:G:O2'	2.07	0.72
16:7I:21:VAL:HG12	16:7I:33:ILE:HD12	1.69	0.72
3:22:57:ILE:HG12	3:22:66:VAL:HG22	1.71	0.72
26:14:395:U:H2'	26:14:396:G:N7	2.03	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:15:42:TRP:O	41:85:64:ARG:NH2	2.19	0.72
1:1G:985:C:N3	1:1G:1220:G:N2	2.37	0.72
26:14:1225:C:O2'	42:95:85:LYS:N	2.23	0.72
1:1G:983:A:N1	1:1G:1222:G:N2	2.38	0.72
26:14:910:A:H62	37:45:12:GLN:HA	1.54	0.72
43:E8:12:ILE:HG13	43:E8:42:ARG:HH11	1.55	0.72
5:4E:45:PHE:CE2	5:4E:47:LYS:HD2	2.25	0.72
9:82:28:VAL:HG22	9:82:63:ILE:HB	1.72	0.72
26:1H:2419:U:H5'	53:O8:23:THR:HG21	1.71	0.72
1:1G:1057:G:H1	1:1G:1203:C:H42	1.37	0.72
28:19:70:TRP:CH2	28:19:150:LYS:HA	2.24	0.72
27:1J:80:U:H2'	27:1J:81:G:N2	2.05	0.72
46:H8:152:ALA:HB1	46:H8:163:LEU:HD11	1.70	0.72
46:H8:165:VAL:HB	46:H8:167:PRO:HD3	1.72	0.72
31:49:67:LYS:H	51:I5:6:HIS:CD2	2.07	0.72
19:AI:5:LEU:HD13	19:AI:10:PHE:HD2	1.54	0.72
38:55:38:VAL:HG12	38:55:42:LYS:HD2	1.71	0.72
39:65:84:GLN:HA	39:65:110:LEU:HD12	1.72	0.72
26:1H:1257:C:H4'	30:31:83:PHE:CD1	2.25	0.72
1:1G:589:C:H42	1:1G:650:G:H1	1.38	0.72
26:1H:141:A:H8	26:1H:1595:G:H21	1.36	0.72
1:1G:532:A:N6	1:1G:1206:G:O2'	2.23	0.72
1:13:1372:U:H5''	9:8E:71:SER:HB2	1.70	0.72
26:14:2134:A:OP2	26:14:2157:G:N2	2.23	0.72
40:B8:3:ARG:HB2	40:B8:6:LEU:HB3	1.70	0.72
36:35:14:LYS:O	36:35:16:ARG:N	2.23	0.72
36:78:50:ARG:HH21	36:78:50:ARG:HG3	1.55	0.71
1:1G:680:C:N4	1:1G:710:G:H1	1.88	0.71
37:88:137:TYR:CE1	46:H8:83:PRO:HG3	2.25	0.71
1:13:1069:C:OP2	58:13:1951:HOH:O	2.08	0.71
46:D5:87:ASP:N	46:D5:87:ASP:OD1	2.21	0.71
26:14:2002:G:N7	58:14:3900:HOH:O	2.22	0.71
26:14:1357:U:O4	58:14:3968:HOH:O	2.03	0.71
26:1H:1386:C:H2'	26:1H:1387:C:C6	2.24	0.71
28:11:146:GLU:HB2	28:11:189:CYS:HB3	1.72	0.71
26:1H:2033:A:H8	58:1H:3948:HOH:O	1.72	0.71
1:13:974:A:OP2	14:5I:41:ARG:NH1	2.23	0.71
26:14:2720:U:H3	26:14:2873:A:H2	1.36	0.71
26:1H:265:A:C8	26:1H:266:G:H1'	2.25	0.71
8:72:120:THR:HG23	8:72:123:GLU:H	1.55	0.71
26:14:2324:C:H5''	26:14:2325:G:H5'	1.70	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:13:1129:C:H4'	1:13:1130:A:H5'	1.72	0.71
24:1L:19:G:N2	24:1L:56:C:N3	2.38	0.71
26:1H:945:A:N3	58:1H:3801:HOH:O	2.24	0.71
1:1G:516:U:O4	58:1G:1725:HOH:O	2.07	0.71
26:1H:1658:C:OP1	58:1H:3694:HOH:O	2.07	0.71
7:62:20:ASP:HB3	7:62:23:VAL:HG23	1.71	0.71
26:14:602:G:O2'	26:14:604:G:O2'	2.06	0.71
32:59:89:ILE:HD12	32:59:130:ARG:HA	1.69	0.71
45:C5:17:SER:HB3	45:C5:71:LYS:HB3	1.71	0.71
45:G8:49:VAL:HG21	45:G8:55:TYR:HE2	1.56	0.71
30:31:32:LEU:HD21	30:31:105:VAL:HG13	1.72	0.71
26:14:619:G:H5''	26:14:620:G:H21	1.56	0.71
1:1G:1305:G:H22	1:1G:1331:G:H2'	1.56	0.71
2:12:178:ARG:HH21	8:72:68:ARG:HH22	1.39	0.71
8:7E:34:GLU:OE1	8:7E:37:ARG:NH1	2.23	0.71
26:1H:958:U:OP2	37:88:14:ARG:NH1	2.23	0.71
10:1I:58:ASP:OD1	58:1I:201:HOH:O	2.09	0.71
26:1H:1772:G:OP1	58:1H:3895:HOH:O	2.08	0.71
37:88:59:ARG:HH11	37:88:59:ARG:H	1.38	0.71
2:12:18:GLY:O	2:12:204:ASN:ND2	2.23	0.71
2:1E:53:ARG:HH12	2:1E:200:ILE:HD12	1.55	0.71
26:1H:1062:G:H2'	26:1H:1063:G:C8	2.26	0.71
26:14:1416:G:O2'	26:14:1417:C:O5'	2.08	0.71
32:51:4:ILE:HD13	32:51:4:ILE:H	1.56	0.71
11:2I:54:ARG:NH1	24:3K:40:C:OP1	2.24	0.71
1:13:835:U:H3	1:13:851:G:H1	1.37	0.71
8:7E:87:SER:HB2	8:7E:93:VAL:HB	1.73	0.71
46:H8:9:TYR:HE1	46:H8:35:ARG:HD3	1.55	0.71
26:14:2439:A:C8	26:14:2439:A:H5'	2.25	0.71
40:B8:77:PRO:HG2	40:B8:80:SER:HB2	1.72	0.71
43:E8:70:TYR:H	43:E8:70:TYR:HD1	1.37	0.71
26:14:270(I):G:H2'	26:14:270(J):G:H8	1.55	0.71
1:1G:853:G:H2'	1:1G:854:G:C8	2.26	0.71
46:H8:9:TYR:CE1	46:H8:35:ARG:HD3	2.26	0.71
26:1H:2298:A:H62	26:1H:2318:G:H8	1.37	0.71
2:1E:178:ARG:HG3	8:7E:72:PRO:HA	1.73	0.71
1:1G:411:A:H62	1:1G:413:G:H21	1.38	0.71
45:G8:82:PRO:HG3	45:G8:97:ARG:HG3	1.73	0.71
26:14:2572:A:C8	29:29:144:ARG:HD2	2.26	0.71
9:8E:112:LYS:HA	9:8E:119:ALA:HB2	1.72	0.71
26:14:1048:A:N6	26:14:1112:G:O2'	2.24	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
45:C5:18:GLY:O	45:C5:21:LYS:N	2.24	0.71
26:14:2127:G:O6	26:14:2161:C:N4	2.18	0.71
1:13:523:A:H61	12:3I:92:ASP:HB2	1.55	0.71
19:AA:39:THR:OG1	19:AA:70:LYS:NZ	2.24	0.71
26:14:2287:A:H62	26:14:2344:U:H3	1.38	0.70
23:2K:54:G:H2'	23:2K:55:5MU:H6	1.54	0.70
26:14:662:G:OP1	36:35:15:ARG:NH2	2.24	0.70
55:M5:53:PRO:O	55:M5:57:ARG:NH1	2.24	0.70
40:B8:50:ILE:HD13	40:B8:64:ARG:HB3	1.71	0.70
26:14:323:G:O2'	26:14:1205:U:N3	2.16	0.70
26:14:479:A:N3	26:14:481:G:H5''	2.06	0.70
24:3K:64:A:N3	24:3K:65:G:N2	2.39	0.70
1:1G:662:G:O2'	1:1G:836:G:OP1	2.10	0.70
26:14:459:U:H5''	54:L5:40:TRP:CD2	2.24	0.70
26:14:831:G:N7	58:14:3657:HOH:O	2.24	0.70
42:95:71:LEU:H	42:95:86:GLY:CA	2.05	0.70
16:7I:26:ARG:HE	16:7I:31:LYS:HB3	1.57	0.70
11:2I:79:SER:HB2	11:2I:106:LYS:HD2	1.73	0.70
36:35:15:ARG:HH21	36:35:17:LYS:HE3	1.57	0.70
26:14:1019:U:OP1	26:14:1035:U:O2'	2.10	0.70
33:61:3:VAL:HG12	33:61:38:LEU:HA	1.73	0.70
28:19:242:ARG:H	28:19:242:ARG:HH11	1.39	0.70
1:13:584:G:O6	58:13:1921:HOH:O	2.06	0.70
19:AA:9:VAL:HG22	51:I5:62:ARG:HB3	1.73	0.70
1:1G:474:G:H2'	1:1G:475:G:C8	2.26	0.70
35:25:2:ILE:HD12	35:25:6:THR:HG21	1.73	0.70
51:M8:12:ALA:HB3	51:M8:24:THR:HB	1.72	0.70
26:1H:2313:C:H4'	31:41:91:ARG:HG3	1.73	0.70
26:1H:1156:A:C8	41:C8:51:LYS:HD3	2.27	0.70
1:13:659:U:H2'	1:13:660:G:H8	1.56	0.70
36:78:121:LYS:HE2	36:78:123:LEU:HD12	1.72	0.70
13:4A:33:ALA:O	13:4A:37:THR:OG1	2.09	0.70
1:13:1348:U:N3	1:13:1374:A:H2	1.88	0.70
1:1G:1368:G:H5'	9:82:112:LYS:HB3	1.73	0.70
1:13:1302:U:OP2	13:4I:21:TYR:OH	2.03	0.70
7:62:67:GLU:OE2	7:62:70:LYS:NZ	2.24	0.70
26:1H:2032:G:H21	29:21:146:THR:CG2	2.03	0.70
26:1H:265:A:H8	26:1H:266:G:H1'	1.56	0.70
32:51:30:LYS:HE3	32:51:81:GLU:H	1.56	0.70
32:59:137:ASP:HB3	32:59:140:LYS:HB2	1.71	0.70
10:1A:51:ARG:HB2	10:1A:60:ARG:HA	1.74	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:1H:1265:A:H3'	52:N8:19:ARG:NH1	2.07	0.70
26:14:943:U:OP2	36:35:36:LYS:HG3	1.92	0.70
36:78:56:SER:HB2	36:78:61:ARG:HD3	1.73	0.70
26:14:2429:G:O6	36:35:61:ARG:NH2	2.25	0.70
30:39:103:LYS:HA	30:39:106:ARG:HG3	1.74	0.70
1:1G:564:C:O2'	8:72:91:ARG:NH2	2.25	0.70
2:1E:15:VAL:HG21	2:1E:209:ARG:HB3	1.73	0.70
38:98:41:ALA:O	38:98:44:LEU:N	2.23	0.70
26:1H:2577:A:H5''	26:1H:2578:G:H5'	1.74	0.69
26:14:1581:G:H2'	26:14:1582:C:O4'	1.92	0.69
30:39:28:ILE:HA	30:39:112:MET:HG2	1.71	0.69
1:13:77:C:N3	1:13:90:C:N4	2.39	0.69
30:31:64:ILE:HG23	30:31:65:TRP:CD1	2.27	0.69
17:8I:67:LYS:HA	17:8I:70:ARG:HH12	1.56	0.69
26:14:1945:G:H2'	26:14:1946:U:C6	2.27	0.69
6:5E:97:PHE:N	18:9I:30:ASP:OD1	2.22	0.69
46:H8:128:VAL:HG23	46:H8:161:VAL:HG11	1.74	0.69
1:13:630:G:H2'	1:13:631:G:O4'	1.92	0.69
26:14:1378:A:O2'	26:14:1380:G:N7	2.23	0.69
26:14:2307:G:O2'	26:14:2308:G:OP2	2.07	0.69
31:49:125:PHE:HB3	31:49:166:ASP:HB2	1.74	0.69
26:14:2327:A:H2'	26:14:2328:A:C8	2.27	0.69
26:14:848:G:H2'	26:14:849:A:H8	1.58	0.69
26:14:607:U:OP1	30:39:102:PRO:HA	1.92	0.69
1:13:1125:U:H5'	1:13:1126:U:H5	1.56	0.69
49:K8:47:ASN:O	49:K8:49:LYS:N	2.24	0.69
26:1H:1021:A:H62	26:1H:1141:U:H3	1.40	0.69
51:M8:13:ARG:O	51:M8:30:GLU:HA	1.92	0.69
36:35:138:LEU:HD12	36:35:144:GLU:HG3	1.74	0.69
26:14:654(E):C:H42	26:14:654(P):G:H22	1.40	0.69
26:14:1329:U:H5''	26:14:1330:C:H5	1.58	0.69
23:2K:47:7MG:H81	23:2K:48:U:H5	1.56	0.69
26:1H:900:A:H3'	26:1H:901:A:H8	1.57	0.69
26:1H:1516:U:H2'	26:1H:1517:G:C8	2.27	0.69
5:4E:126:ARG:HH11	5:4E:126:ARG:HG3	1.57	0.69
26:1H:1678:G:H21	26:1H:1989:G:H22	1.37	0.69
26:1H:76:C:O2'	49:K8:62:THR:HG21	1.92	0.69
40:75:4:GLY:O	40:75:7:ILE:N	2.25	0.69
45:C5:77:PRO:HD2	45:C5:103:GLY:HA2	1.73	0.69
26:1H:1676:A:OP2	58:1H:3703:HOH:O	2.09	0.69
26:1H:573:G:O2'	26:1H:574:C:H3'	1.92	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
37:88:135:ASP:HB3	37:88:137:TYR:H	1.56	0.69
26:14:987:G:O2'	26:14:1000:A:N3	2.21	0.69
15:6A:33:THR:HG21	15:6A:85:LEU:HD22	1.72	0.69
26:1H:2683:C:OP1	40:B8:53:ARG:NH2	2.26	0.69
28:19:40:THR:OG1	28:19:41:GLY:N	2.25	0.69
12:3I:7:ILE:HD13	12:3I:10:LEU:HD12	1.73	0.69
26:1H:1113:U:H5'	32:51:2:SER:HB2	1.72	0.69
37:88:82:ARG:HD2	37:88:82:ARG:N	2.08	0.69
26:1H:2308:G:N1	26:1H:2311:A:H2	1.84	0.69
26:14:1022:G:O2'	26:14:1023:U:OP2	2.05	0.69
39:A8:89:ARG:HG2	39:A8:89:ARG:O	1.92	0.69
26:14:2685:G:C6	58:14:3783:HOH:O	2.46	0.69
26:1H:1174:A:H1'	26:1H:1178:C:H41	1.57	0.69
9:8E:13:ALA:HB2	9:8E:68:GLY:HA3	1.75	0.69
32:59:106:THR:HG22	32:59:112:PRO:HB3	1.75	0.69
29:29:33:VAL:HG23	29:29:47:VAL:HG13	1.73	0.69
26:14:882:G:H22	26:14:894:C:H42	1.40	0.69
1:1G:1014:A:H4'	19:AA:14:HIS:CE1	2.27	0.69
1:13:963:G:N3	10:1I:55:LYS:NZ	2.40	0.69
13:4I:81:LEU:HD22	13:4I:88:ARG:HB2	1.72	0.69
19:AI:5:LEU:HD13	19:AI:10:PHE:CD2	2.27	0.69
12:3A:117:ARG:HB3	12:3A:122:THR:HB	1.74	0.69
31:41:131:TYR:O	31:41:159:VAL:HG22	1.93	0.69
1:1G:1376:U:H2'	1:1G:1377:A:H8	1.56	0.69
26:1H:2134:A:OP2	26:1H:2157:G:N2	2.25	0.69
12:3A:41:ARG:HH11	12:3A:41:ARG:HB3	1.58	0.69
1:1G:943:U:H1'	9:82:124:GLN:HE22	1.57	0.69
26:14:1676:A:OP2	58:14:3549:HOH:O	2.11	0.69
4:3E:26:CYS:HA	4:3E:31:CYS:HB2	1.74	0.69
27:1J:42:C:O2'	31:49:67:LYS:O	2.10	0.69
44:F8:51:VAL:HG13	44:F8:81:VAL:HG23	1.75	0.69
26:14:1689:A:H62	26:14:1698:A:H2	1.40	0.69
7:62:21:VAL:HG23	7:62:22:LEU:HD12	1.74	0.69
39:A8:11:LYS:HD3	39:A8:91:PRO:HD3	1.75	0.69
26:1H:863:A:N7	58:1H:3909:HOH:O	2.25	0.69
34:58:96:GLU:O	34:58:98:VAL:N	2.26	0.69
26:1H:2136:C:N3	26:1H:2155:G:N1	2.38	0.69
1:1G:1256:A:N6	1:1G:1278:U:OP2	2.24	0.69
26:14:2025:C:N4	58:14:3843:HOH:O	2.26	0.69
17:8I:22:LEU:HD11	17:8I:39:SER:HB3	1.74	0.69
26:1H:1603:A:OP1	58:1H:3648:HOH:O	2.09	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
10:1I:27:ALA:HB1	10:1I:34:VAL:HG11	1.74	0.68
40:75:8:LYS:HZ2	40:75:8:LYS:HB2	1.58	0.68
3:2E:40:ARG:O	3:2E:44:GLU:HG2	1.92	0.68
49:G5:47:ASN:O	49:G5:49:LYS:N	2.24	0.68
40:75:50:ILE:HD11	40:75:102:ILE:HD11	1.73	0.68
1:13:1453:G:H2'	20:BI:39:LYS:HE2	1.74	0.68
51:I5:38:LYS:HA	51:I5:44:THR:HG21	1.75	0.68
26:1H:242:G:OP1	58:1H:4111:HOH:O	2.10	0.68
26:1H:386:G:OP2	58:1H:4346:HOH:O	2.11	0.68
55:Q8:23:VAL:HG13	55:Q8:46:ARG:HG3	1.75	0.68
26:14:881:G:O6	26:14:882:G:N2	2.26	0.68
28:11:223:GLY:HA3	28:11:231:HIS:ND1	2.08	0.68
1:13:1001:G:N2	1:13:1040:U:O2	2.27	0.68
30:31:12:LEU:O	30:31:127:GLU:N	2.25	0.68
7:6E:20:ASP:HB3	7:6E:23:VAL:HG23	1.75	0.68
26:1H:1786:A:H2	26:1H:2606:C:H1'	1.59	0.68
39:A8:35:ILE:HG22	39:A8:97:ARG:HH21	1.59	0.68
1:1G:1352:C:H42	1:1G:1370:G:H1	1.42	0.68
1:1G:501:C:H2'	1:1G:502:G:H8	1.58	0.68
8:7E:77:GLU:HG2	8:7E:78:GLN:H	1.59	0.68
28:11:242:ARG:O	58:11:401:HOH:O	2.11	0.68
51:I5:20:ASN:ND2	51:I5:39:CYS:SG	2.66	0.68
50:L8:56:VAL:HG12	50:L8:57:GLU:H	1.58	0.68
46:D5:111:VAL:HG22	46:D5:112:ARG:HG2	1.74	0.68
26:14:2632:A:HO2'	26:14:2811:G:HO2'	1.21	0.68
26:1H:1189:A:OP2	58:1H:3820:HOH:O	2.11	0.68
36:35:55:ARG:HG2	36:35:56:SER:H	1.57	0.68
30:39:102:PRO:HB2	30:39:105:VAL:HG23	1.74	0.68
23:2K:47:7MG:H81	23:2K:48:U:C5	2.29	0.68
1:1G:1014:A:H2'	1:1G:1015:A:C8	2.29	0.68
26:1H:2096:U:H3	26:1H:2193:G:H1	1.42	0.68
26:1H:357:A:H2'	26:1H:358:U:H6	1.58	0.68
40:B8:56:GLY:O	40:B8:59:THR:HG22	1.94	0.68
10:1A:48:THR:HA	10:1A:62:HIS:HB3	1.76	0.68
26:14:1332:G:H5'	26:14:1332:G:C8	2.29	0.68
19:AA:18:LYS:HA	19:AA:21:GLU:HG2	1.74	0.68
46:D5:163:LEU:HD23	46:D5:163:LEU:H	1.58	0.68
26:1H:2136:C:N4	26:1H:2155:G:O6	2.20	0.68
1:13:1391:U:H2'	1:13:1392:G:C8	2.29	0.68
26:14:71:A:H2	44:B5:31:HIS:NE2	1.90	0.68
55:Q8:7:HIS:HB3	55:Q8:58:ILE:HG23	1.76	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:1H:192:C:OP1	58:1H:4127:HOH:O	2.11	0.68
26:14:274:G:H2'	26:14:275:G:H4'	1.76	0.68
1:13:101:A:H2'	1:13:102:G:H8	1.58	0.68
26:14:2777:G:H5''	26:14:2778:A:H5'	1.76	0.68
1:1G:1238:A:OP1	1:1G:1335:C:O2'	2.12	0.68
1:13:1319:A:O2'	1:13:1323:G:N7	2.22	0.68
26:1H:578:A:OP2	58:1H:4150:HOH:O	2.11	0.68
44:F8:5:TYR:O	49:K8:36:ARG:NH2	2.27	0.68
26:1H:1664:A:OP2	58:1H:3962:HOH:O	2.10	0.68
26:1H:2057:A:OP2	58:1H:3617:HOH:O	2.12	0.68
36:78:26:GLY:O	58:78:302:HOH:O	2.12	0.68
19:AA:60:VAL:HG21	19:AA:74:PHE:HB3	1.73	0.68
2:1E:87:ARG:NH2	2:1E:220:ASP:OD1	2.27	0.68
11:2A:29:ILE:HG22	11:2A:44:SER:HB2	1.76	0.68
48:F5:20:ARG:HG2	48:F5:20:ARG:HH11	1.58	0.68
3:2E:15:THR:HG22	3:2E:16:ARG:H	1.59	0.68
26:1H:1163:G:H2'	26:1H:1164:G:H8	1.59	0.68
26:1H:1697:G:OP2	26:1H:1698:A:O2'	2.11	0.68
1:13:1454:G:OP1	20:BI:39:LYS:NZ	2.18	0.68
44:F8:5:TYR:CE1	49:K8:30:ARG:HG3	2.29	0.68
30:31:167:ALA:HB1	30:31:173:VAL:HG11	1.74	0.68
26:1H:2695:C:H2'	26:1H:2696:U:H6	1.59	0.68
1:1G:1513:A:H2'	1:1G:1514:C:C6	2.28	0.68
26:1H:2685:G:O6	58:1H:4249:HOH:O	2.08	0.68
26:14:2393:A:H4'	36:35:62:LEU:H	1.58	0.68
13:4I:94:ARG:HH22	26:1H:887:A:H5''	1.56	0.68
26:14:2017:U:OP1	58:14:3835:HOH:O	2.12	0.68
23:2L:24:C:H2'	23:2L:25:U:C6	2.29	0.68
1:13:1178:G:OP2	9:8E:93:ARG:NH2	2.26	0.68
38:98:83:ILE:HG22	38:98:87:TYR:HE2	1.58	0.68
29:29:111:ARG:HA	38:55:2:ARG:HH12	1.59	0.68
26:14:2074:U:OP1	58:14:3509:HOH:O	2.11	0.68
34:58:96:GLU:C	34:58:98:VAL:H	1.98	0.68
26:1H:270(V):G:H2'	26:1H:270(W):G:H8	1.59	0.68
26:1H:2709:G:O2'	58:1H:3672:HOH:O	2.12	0.68
1:13:1145:C:H4'	1:13:1146:A:H8	1.58	0.68
1:1G:1190:G:H5'	3:22:176:HIS:CE1	2.29	0.68
1:1G:516:U:O4	58:1G:1729:HOH:O	2.09	0.67
3:2E:19:GLU:O	3:2E:40:ARG:NH2	2.27	0.67
44:F8:61:GLY:N	44:F8:75:ASP:OD1	2.27	0.67
1:13:128:G:H5'	17:8I:2:PRO:O	1.93	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
15:6I:74:ASP:HB3	15:6I:77:ARG:HG2	1.76	0.67
29:21:23:VAL:HA	29:21:185:LYS:HA	1.76	0.67
5:42:81:GLU:HB3	5:42:90:VAL:HG13	1.75	0.67
26:1H:1329:U:H5''	26:1H:1330:C:H5	1.59	0.67
15:6A:87:ILE:HG22	15:6A:88:ARG:H	1.59	0.67
55:Q8:49:VAL:HG11	55:Q8:52:LYS:HG3	1.77	0.67
26:14:1060:U:H4'	26:14:1061:U:H5''	1.77	0.67
43:A5:72:LYS:HB3	43:A5:106:ILE:HG13	1.76	0.67
1:1G:837:G:H1	1:1G:849:C:H42	1.42	0.67
2:1E:73:THR:O	2:1E:78:GLN:NE2	2.27	0.67
26:14:469:G:O6	54:L5:37:LYS:HE2	1.94	0.67
26:1H:1406:U:H2'	26:1H:1407:C:C6	2.28	0.67
45:C5:68:HIS:HB3	45:C5:71:LYS:HG3	1.77	0.67
5:4E:8:GLU:OE1	5:4E:63:ARG:NH2	2.27	0.67
45:G8:76:CYS:SG	45:G8:97:ARG:HG2	2.35	0.67
26:1H:2815:C:H5'	52:N8:29:THR:HG21	1.76	0.67
26:1H:593:G:O3'	55:Q8:60:LEU:HD22	1.95	0.67
26:14:1614:A:OP1	58:14:3514:HOH:O	2.13	0.67
16:7I:51:VAL:HG11	16:7I:77:ALA:HB1	1.77	0.67
23:2L:24:C:H2'	23:2L:25:U:H6	1.58	0.67
1:13:1133:G:H22	1:13:1141:C:H42	1.41	0.67
24:3L:5:G:H22	24:3L:68:C:H42	1.41	0.67
26:1H:2212:A:H1'	26:1H:2215:G:C5	2.30	0.67
41:85:90:VAL:HG22	42:95:39:LEU:HB3	1.74	0.67
28:11:10:THR:OG1	28:11:13:ARG:HB2	1.94	0.67
40:B8:111:ARG:H	40:B8:111:ARG:HD3	1.58	0.67
42:95:85:LYS:HG3	42:95:87:HIS:N	2.09	0.67
26:1H:192:C:OP1	58:1H:4130:HOH:O	2.12	0.67
2:12:163:PHE:HD1	2:12:185:ILE:HG13	1.60	0.67
26:1H:459:U:H5''	54:P8:40:TRP:CD2	2.29	0.67
1:13:129(A):G:H4'	1:13:130:A:H5''	1.77	0.67
26:14:2294:C:OP2	39:65:89:ARG:NH2	2.28	0.67
5:42:61:TYR:HA	5:42:64:ARG:HB2	1.76	0.67
26:1H:1194:A:OP2	26:1H:1194:A:H8	1.78	0.67
26:14:71:A:C8	26:14:71:A:H5'	2.29	0.67
2:12:42:ILE:HD11	2:12:202:PRO:HB2	1.77	0.67
26:14:892:G:N2	26:14:894:C:OP1	2.28	0.67
43:A5:73:ALA:HB3	43:A5:106:ILE:HD11	1.75	0.67
13:4A:13:LYS:HA	13:4A:44:ARG:HH11	1.60	0.67
38:55:67:LEU:HD12	38:55:76:VAL:HG21	1.75	0.67
26:1H:2346:A:O2'	53:O8:24:GLU:OE2	2.13	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
13:4A:3:ARG:HB2	51:I5:34:GLU:HG3	1.76	0.67
38:98:97:VAL:HG22	38:98:114:VAL:HG22	1.75	0.67
26:14:661:C:H1'	36:35:12:ALA:HA	1.76	0.67
26:14:198:C:H5'	26:14:2244:U:OP1	1.94	0.67
42:95:48:GLY:N	42:95:52:VAL:HG22	2.10	0.67
1:13:524:G:H2'	1:13:525:C:C6	2.30	0.67
29:21:18:ASP:HB3	40:B8:82:LEU:HD11	1.76	0.67
5:4E:110:LEU:HD13	5:4E:118:ILE:HD13	1.77	0.67
28:19:255:LYS:CE	28:19:255:LYS:H	2.08	0.67
38:55:74:LYS:HE2	38:55:77:ARG:HH21	1.59	0.67
26:14:993:G:N3	42:95:89:GLN:NE2	2.38	0.67
46:D5:17:ALA:HA	46:D5:20:ARG:HD2	1.75	0.67
29:21:116:VAL:O	29:21:117:MET:HB3	1.94	0.67
26:1H:2271:G:N7	58:1H:4058:HOH:O	2.27	0.67
40:B8:20:PRO:HD2	40:B8:86:ILE:HG23	1.77	0.67
26:1H:860:U:C5	26:1H:917:A:C2	2.82	0.67
37:45:27:VAL:HG11	46:D5:81:ARG:NH2	2.09	0.67
26:1H:662:G:H5'	36:78:15:ARG:H	1.60	0.67
37:45:22:LYS:HB3	37:45:23:GLY:HA2	1.76	0.67
46:D5:44:PHE:HE1	46:D5:88:PHE:HZ	1.43	0.67
37:88:106:VAL:HG21	37:88:114:ALA:HB1	1.76	0.67
1:13:1435:G:H2'	1:13:1436:U:C6	2.30	0.67
37:45:57:HIS:NE2	37:45:116:GLU:HB3	2.09	0.67
26:1H:981:A:OP1	58:1H:4157:HOH:O	2.12	0.67
31:41:77:ILE:HG22	31:41:82:LEU:HD12	1.77	0.67
26:14:1485:G:H1	26:14:1504:C:N4	1.93	0.67
26:1H:450:G:OP2	58:1H:3833:HOH:O	2.12	0.67
1:1G:474:G:H2'	1:1G:475:G:H8	1.59	0.67
1:1G:458:C:N3	1:1G:474:G:N2	2.43	0.67
40:75:1:MET:HA	40:75:5:ALA:HB3	1.76	0.67
27:16:8:U:O2	27:16:112:G:N1	2.19	0.67
26:14:630:G:N2	26:14:633:A:OP2	2.28	0.67
30:39:4:VAL:HA	30:39:19:GLU:HB3	1.77	0.67
26:14:453:C:OP1	58:14:3864:HOH:O	2.13	0.67
25:4L:13:A:O2'	25:4L:14:A:OP1	2.13	0.67
42:95:35:LEU:O	42:95:37:VAL:HG22	1.94	0.67
1:1G:456:C:N3	1:1G:476:G:N2	2.38	0.67
26:1H:1676:A:OP2	58:1H:3698:HOH:O	2.12	0.67
1:1G:644:G:H4'	8:72:92:ARG:HH12	1.60	0.67
26:1H:912:C:OP1	37:88:8:LYS:NZ	2.25	0.67
46:H8:147:GLY:N	46:H8:174:VAL:O	2.25	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1G:1177:G:OP2	9:82:97:LYS:NZ	2.27	0.67
34:15:15:LEU:HB2	34:15:134:ARG:HG2	1.76	0.67
9:82:10:ARG:NH1	9:82:105:ASP:OD2	2.28	0.67
1:13:1023:G:H3'	1:13:1024:G:H5''	1.77	0.66
26:1H:2057:A:OP2	58:1H:3613:HOH:O	2.13	0.66
1:1G:1286:A:H5'	21:1B:25:LYS:HD3	1.77	0.66
26:1H:1839:G:OP2	58:1H:4115:HOH:O	2.12	0.66
26:1H:606:U:OP2	30:31:104:LYS:NZ	2.29	0.66
26:1H:2164:C:OP2	26:1H:2166:G:N2	2.29	0.66
4:32:59:ARG:O	4:32:63:LYS:N	2.27	0.66
26:14:34:C:O2'	26:14:35:G:OP1	2.12	0.66
26:1H:31:C:OP1	58:1H:3738:HOH:O	2.12	0.66
3:22:81:GLY:HA2	3:22:85:ARG:HD3	1.76	0.66
13:4A:86:CYS:HB2	19:AA:73:GLU:HB3	1.76	0.66
39:A8:24:LEU:HD12	39:A8:41:ASP:HB2	1.75	0.66
26:1H:308:G:N7	58:1H:4397:HOH:O	2.27	0.66
26:14:2273:A:H2'	26:14:2274:A:C8	2.31	0.66
26:1H:1061:U:H4'	26:1H:1070:A:H1'	1.77	0.66
1:1G:1081:G:OP1	5:42:18:ARG:HB2	1.95	0.66
26:14:1226:G:OP1	42:95:69:LYS:NZ	2.17	0.66
1:13:859:A:H2'	1:13:860:A:C8	2.31	0.66
26:1H:1140:C:OP1	34:58:23:LEU:HB3	1.94	0.66
1:1G:1235:U:O2'	1:1G:1305:G:O5'	2.13	0.66
13:4I:23:TYR:CE2	13:4I:71:ARG:HG3	2.30	0.66
1:1G:1173:G:OP1	7:62:5:ARG:NH2	2.28	0.66
26:14:21:A:H61	26:14:519:U:H3	1.41	0.66
31:41:21:ARG:HG2	31:41:21:ARG:HH11	1.59	0.66
14:5I:24:CYS:HB3	14:5I:27:CYS:HB2	1.76	0.66
1:1G:1274:G:H2'	1:1G:1275:A:H8	1.60	0.66
28:19:2:ALA:HB3	28:19:20:ASP:HB2	1.76	0.66
26:1H:1514:U:H2'	26:1H:1515:C:H6	1.60	0.66
1:1G:957:U:O2'	1:1G:959:A:N7	2.26	0.66
26:1H:1405:U:H2'	26:1H:1406:U:C6	2.30	0.66
26:14:2701:C:H3'	26:14:2702:U:C5'	2.24	0.66
27:1J:14:U:O2'	27:1J:107:U:O2'	2.14	0.66
4:32:199:ASN:HB3	4:32:202:LEU:HG	1.77	0.66
1:1G:1028:C:H42	1:1G:1033:G:H1	1.44	0.66
26:14:1839:G:OP2	58:14:3688:HOH:O	2.12	0.66
30:39:79:GLY:HA2	30:39:86:GLY:HA2	1.77	0.66
1:13:1297:C:OP1	13:4I:13:LYS:NZ	2.28	0.66
26:14:2130:U:H2'	26:14:2158:A:N1	2.11	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:6E:16:LEU:HD11	9:8E:45:ALA:HB2	1.76	0.66
13:4I:94:ARG:NH1	26:1H:888:C:OP1	2.25	0.66
1:13:396:G:O2'	1:13:398:C:OP1	2.08	0.66
26:1H:2306:C:H3'	26:1H:2307:G:H5'	1.77	0.66
22:1K:18:G:H4'	22:1K:19:G:OP1	1.96	0.66
26:1H:86:C:H4'	26:1H:104:U:H1'	1.76	0.66
26:1H:2058:A:N7	58:1H:3615:HOH:O	2.27	0.66
30:39:167:ALA:HB1	30:39:173:VAL:HG11	1.77	0.66
31:49:20:ILE:HA	31:49:25:TYR:HB2	1.77	0.66
26:14:1537:C:O2'	26:14:1538:G:O4'	2.13	0.66
1:1G:742:G:OP2	15:6A:35:ARG:NH2	2.29	0.66
26:1H:229:A:H4'	26:1H:230:U:H5'	1.76	0.66
1:13:991:U:O4	1:13:1212:U:O2'	2.12	0.66
1:1G:999:U:H2'	1:1G:1000:A:C8	2.30	0.66
26:14:1639:U:OP1	58:14:3555:HOH:O	2.13	0.66
26:1H:571:A:OP2	58:1H:3813:HOH:O	2.14	0.66
26:1H:1800:C:OP1	28:11:266:SER:OG	2.11	0.66
26:14:1019:U:H2'	26:14:1020:A:C8	2.31	0.66
26:1H:2502:G:OP2	58:1H:3632:HOH:O	2.14	0.66
1:13:963:G:H1	1:13:972:C:H42	1.41	0.66
1:1G:1213:A:N6	1:1G:1215:G:N3	2.43	0.66
32:51:4:ILE:HG13	32:51:6:ARG:NE	2.11	0.66
1:1G:411:A:C5	1:1G:413:G:H1'	2.30	0.66
26:14:249:C:OP1	58:14:3522:HOH:O	2.13	0.66
26:14:2537:U:H2'	26:14:2538:C:C6	2.30	0.66
1:13:1062:U:H2'	1:13:1063:C:C6	2.31	0.66
26:14:1428:C:N4	26:14:1570:A:OP2	2.23	0.66
28:11:96:HIS:CE1	28:11:102:LYS:HE2	2.31	0.66
29:29:55:ASN:O	29:29:57:LYS:NZ	2.29	0.66
46:H8:165:VAL:HB	46:H8:166:SER:HA	1.77	0.66
1:13:972:C:OP1	58:13:1881:HOH:O	2.13	0.66
26:14:273(C):C:N4	26:14:363(C):G:H1	1.94	0.66
30:39:123:LEU:O	30:39:125:LEU:N	2.25	0.66
29:29:102:VAL:HG12	29:29:200:GLU:HA	1.77	0.66
31:41:173:LEU:HB3	31:41:178:PHE:HD2	1.61	0.66
9:82:4:TYR:HB2	9:82:19:LEU:HD12	1.78	0.66
38:98:12:ARG:HD3	38:98:16:HIS:CD2	2.30	0.66
23:2L:41:C:H2'	23:2L:42:C:H6	1.61	0.66
33:61:132:PRO:O	33:61:133:HIS:ND1	2.29	0.66
51:I5:16:CYS:HA	51:I5:33:VAL:HG13	1.77	0.66
1:13:736:C:H2'	1:13:737:A:C8	2.31	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:252:G:OP2	36:35:50:ARG:NH2	2.29	0.65
1:13:1034:G:N2	1:13:1035:A:N7	2.43	0.65
1:13:235:C:H5'	17:8I:70:ARG:HG2	1.77	0.65
2:1E:111:ARG:HG2	2:1E:111:ARG:HH11	1.60	0.65
38:98:67:LEU:HD22	38:98:76:VAL:HG21	1.78	0.65
40:75:24:PRO:HD3	40:75:52:ILE:HD12	1.78	0.65
4:32:64:LEU:HB2	4:32:198:VAL:HG11	1.78	0.65
26:14:330:A:H2	26:14:1210:A:HO2'	1.44	0.65
1:13:272:C:H2'	1:13:273:A:H8	1.60	0.65
26:14:1225:C:H4'	42:95:85:LYS:HB2	1.78	0.65
42:95:85:LYS:HD2	42:95:86:GLY:H	1.61	0.65
26:1H:1658:C:OP1	58:1H:3695:HOH:O	2.14	0.65
31:49:11:TYR:OH	31:49:16:ARG:NH2	2.28	0.65
26:14:1341:U:OP2	26:14:1394:U:O2'	2.12	0.65
9:82:27:THR:HB	9:82:32:ASP:HA	1.79	0.65
26:1H:2781:A:H5''	26:1H:2782:G:H5'	1.77	0.65
26:14:309:G:N3	26:14:329:G:O2'	2.30	0.65
10:1A:12:ASP:OD1	10:1A:13:HIS:N	2.29	0.65
1:1G:661:G:H1	1:1G:744:C:H42	1.43	0.65
13:4A:94:ARG:CZ	19:AA:78:ARG:HH22	2.09	0.65
26:1H:507:A:H5''	26:1H:508:G:H5'	1.78	0.65
5:42:88:LYS:HB3	5:42:123:LEU:HB2	1.78	0.65
26:14:2845:G:N2	26:14:2871:C:O2	2.27	0.65
26:1H:2593:U:H2'	26:1H:2594:C:C6	2.31	0.65
28:19:69:ARG:NH2	28:19:128:GLY:O	2.29	0.65
7:6E:15:ASP:HB3	7:6E:20:ASP:H	1.61	0.65
26:1H:2125:G:N2	26:1H:2172:U:OP1	2.27	0.65
1:1G:947:G:O3'	13:4A:109:THR:OG1	2.13	0.65
46:H8:45:ASP:OD2	46:H8:49:ARG:NH1	2.29	0.65
30:39:130:ALA:H	30:39:142:TRP:HD1	1.42	0.65
38:55:29:LEU:HB3	38:55:75:LEU:HD21	1.78	0.65
26:1H:2061:G:H5'	58:1H:3632:HOH:O	1.96	0.65
26:14:2331:G:H4'	47:E5:43:THR:H	1.61	0.65
32:51:6:ARG:HB3	32:51:65:HIS:CG	2.31	0.65
1:1G:683:G:N2	1:1G:707:C:O2	2.30	0.65
4:32:61:LYS:HE2	4:32:206:PHE:CE2	2.32	0.65
26:1H:2636:U:OP1	29:21:79:ARG:HA	1.97	0.65
1:13:1263:C:H2'	1:13:1264:C:H6	1.61	0.65
26:1H:2580:U:H4'	29:21:130:GLY:HA3	1.78	0.65
26:1H:287:C:H2'	26:1H:288:C:H6	1.62	0.65
26:1H:2327:A:H2'	26:1H:2328:A:C8	2.30	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
27:1J:70:C:H2'	27:1J:71:C:H6	1.61	0.65
26:14:1188:U:O2'	26:14:1189:A:H5'	1.96	0.65
26:1H:1438:U:H2'	26:1H:1439:A:H8	1.61	0.65
26:1H:607:U:OP1	30:31:102:PRO:HA	1.97	0.65
26:1H:2502:G:OP2	58:1H:3633:HOH:O	2.13	0.65
26:1H:1141:U:H6	34:58:63:THR:HG1	1.44	0.65
26:1H:1332:G:N2	26:1H:1609:A:O2'	2.30	0.65
19:AI:50:ALA:HB1	19:AI:57:HIS:HB3	1.77	0.65
27:1J:6:C:H2'	27:1J:7:G:H5''	1.78	0.65
39:A8:59:LYS:HG2	39:A8:60:GLY:H	1.60	0.65
47:I8:53:MET:HG3	47:I8:59:LEU:CD2	2.26	0.65
48:F5:89:GLU:HA	48:F5:93:GLU:HG3	1.77	0.65
26:1H:226:G:H21	26:1H:228:A:H2	1.45	0.65
26:1H:1342:A:OP2	58:1H:4224:HOH:O	2.14	0.65
1:13:1177:G:OP1	1:13:1177:G:H4'	1.95	0.65
49:K8:47:ASN:HB2	49:K8:50:ILE:HD11	1.79	0.65
26:1H:2115:G:N2	26:1H:2172:U:O2	2.30	0.65
27:16:8:U:N3	27:16:112:G:O6	2.19	0.65
2:1E:5:ILE:HG13	2:1E:6:THR:HG22	1.78	0.65
1:13:448:A:OP2	1:13:485:G:N2	2.23	0.65
26:14:1670:C:O2	29:29:129:HIS:NE2	2.29	0.65
26:14:599:G:N7	58:14:4021:HOH:O	2.29	0.65
26:1H:2232:U:P	48:J8:40:ARG:HH12	2.20	0.65
12:3I:24:VAL:HB	12:3I:27:LEU:HD12	1.79	0.65
33:61:92:VAL:HG13	33:61:120:ILE:HG23	1.78	0.65
26:1H:2285:C:OP2	53:O8:28:ARG:HG3	1.96	0.65
8:72:110:ALA:H	8:72:121:ASP:HB3	1.61	0.65
29:29:47:VAL:HG21	29:29:86:PRO:HD2	1.78	0.65
26:14:2645:G:H3'	26:14:2646:C:H5'	1.78	0.65
47:I8:11:ARG:O	47:I8:14:ARG:NH2	2.29	0.65
45:G8:53:PRO:HA	45:G8:56:PRO:HG3	1.79	0.65
26:14:2256:G:N7	58:14:3650:HOH:O	2.29	0.65
7:62:113:GLU:HB2	7:62:119:ARG:HG2	1.77	0.65
26:1H:2070:G:OP1	58:1H:4121:HOH:O	2.14	0.65
50:H5:40:THR:HG23	50:H5:43:ILE:HG12	1.77	0.65
4:32:23:GLY:H	4:32:26:CYS:HB2	1.62	0.65
36:35:55:ARG:HG2	36:35:56:SER:N	2.12	0.65
26:14:2738:A:OP2	58:14:3999:HOH:O	2.14	0.65
26:1H:624:C:OP1	58:1H:4113:HOH:O	2.14	0.65
26:1H:1447:G:N7	58:1H:4360:HOH:O	2.30	0.65
28:11:136:ILE:O	28:11:168:ARG:NH2	2.30	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:52:70:ASP:OD1	6:52:70:ASP:N	2.30	0.65
26:14:2745:C:O2	32:59:139:GLN:NE2	2.24	0.65
27:16:7:G:H4'	39:A8:29:PHE:CD2	2.31	0.65
39:65:14:VAL:HG11	39:65:89:ARG:HE	1.62	0.65
42:95:43:GLU:O	42:95:45:THR:OG1	2.15	0.65
28:11:59:LYS:HD2	28:11:60:ARG:H	1.61	0.65
26:14:629:G:H1	26:14:634:C:H42	1.43	0.65
1:1G:538:G:H5''	12:3A:114:LYS:HB2	1.78	0.65
1:1G:579:G:O3'	15:6A:54:ARG:NH2	2.30	0.64
31:41:67:LYS:HE2	51:M8:6:HIS:CE1	2.32	0.64
1:13:659:U:H2'	1:13:660:G:C8	2.32	0.64
1:1G:426:G:OP1	4:32:36:ARG:NH2	2.30	0.64
26:14:34:C:HO2'	26:14:35:G:P	2.20	0.64
7:6E:91:VAL:HB	7:6E:96:GLN:HG2	1.79	0.64
26:14:96:G:H4'	49:G5:48:HIS:CE1	2.33	0.64
36:35:85:LEU:HA	36:35:88:LEU:HB3	1.79	0.64
26:1H:1441:G:H2'	26:1H:1442:G:C8	2.32	0.64
30:31:66:PRO:O	30:31:67:GLN:HB3	1.96	0.64
1:1G:1342:C:H4'	9:82:125:TYR:HB3	1.78	0.64
26:1H:1332:G:H21	26:1H:1610:A:H8	1.45	0.64
8:7E:21:LYS:O	8:7E:65:TYR:OH	2.14	0.64
1:13:1224:G:C6	1:13:1322:C:H1'	2.32	0.64
1:13:262:A:H2'	1:13:263:A:C8	2.32	0.64
26:1H:1024:G:H3'	26:1H:1025:G:H5''	1.79	0.64
40:B8:58:ASN:ND2	40:B8:58:ASN:O	2.26	0.64
26:1H:1036:G:H1	26:1H:1119:C:H42	1.46	0.64
26:14:2689:U:OP2	26:14:2719:G:N2	2.31	0.64
1:1G:1348:U:H4'	9:82:120:ARG:HD2	1.77	0.64
29:21:77:ILE:O	29:21:79:ARG:N	2.29	0.64
26:1H:570:G:OP2	58:1H:3811:HOH:O	2.14	0.64
11:2I:32:ILE:HD11	11:2I:68:ALA:HB1	1.79	0.64
1:1G:406:G:H1	1:1G:436:C:H42	1.46	0.64
26:14:1327:C:OP2	58:14:3607:HOH:O	2.15	0.64
55:Q8:52:LYS:HA	55:Q8:54:GLU:HB2	1.79	0.64
52:J5:16:ARG:NH1	52:J5:17:ASP:OD1	2.30	0.64
26:14:635:C:O2'	26:14:639:U:OP1	2.15	0.64
5:42:6:PHE:HB2	5:42:34:VAL:HG22	1.80	0.64
26:14:2272:U:O4	58:14:3727:HOH:O	2.09	0.64
2:12:190:THR:O	2:12:191:ASP:HB3	1.98	0.64
29:29:66:HIS:HB3	29:29:68:ALA:HB2	1.79	0.64
26:14:1945:G:H2'	26:14:1946:U:H6	1.62	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:1H:2286:A:N6	53:O8:24:GLU:OE1	2.30	0.64
32:51:15:VAL:HG12	32:51:29:PRO:HD2	1.79	0.64
1:13:1291:G:OP1	7:6E:37:ASN:ND2	2.30	0.64
41:85:91:ASP:OD1	41:85:96:ALA:HB2	1.97	0.64
26:14:2415:G:H4'	36:35:67:MET:N	2.13	0.64
1:1G:1095:U:P	1:1G:1108:G:H1	2.20	0.64
1:13:390:C:H2'	1:13:391:G:C8	2.33	0.64
1:13:405:U:O4	4:3E:2:GLY:N	2.31	0.64
1:13:412:A:H4'	1:13:413:G:O5'	1.98	0.64
31:49:50:ALA:HB2	31:49:87:PRO:HG3	1.79	0.64
2:1E:17:PHE:HB3	2:1E:44:LEU:HD21	1.79	0.64
29:29:11:MET:HG3	29:29:24:THR:H	1.62	0.64
17:8I:76:LEU:HD11	17:8I:79:SER:HB3	1.79	0.64
29:29:29:GLY:H	29:29:51:PHE:HE1	1.43	0.64
26:14:731:C:OP1	58:14:3866:HOH:O	2.15	0.64
26:1H:270(E):G:H1	26:1H:270(U):C:H42	1.46	0.64
26:1H:2839:G:H21	38:98:92:GLY:HA3	1.62	0.64
35:25:104:ARG:HH12	40:75:36:GLU:HB3	1.63	0.64
28:11:182:LEU:H	28:11:272:ALA:HB3	1.62	0.64
38:98:51:LEU:HD22	38:98:66:VAL:HG13	1.79	0.64
1:13:828:A:H2'	1:13:829:G:O4'	1.97	0.64
26:14:1019:U:H2'	26:14:1020:A:H8	1.62	0.64
26:14:994:C:OP1	41:85:53:ARG:NH2	2.30	0.64
42:95:70:ILE:N	42:95:86:GLY:O	2.19	0.64
27:16:43:C:H5''	51:M8:1:MET:HG2	1.79	0.64
17:8I:66:SER:O	17:8I:70:ARG:NH1	2.31	0.64
1:1G:1295:G:O2'	13:4A:14:ARG:NH1	2.30	0.64
2:12:84:GLU:HB3	2:12:219:VAL:HG11	1.79	0.64
42:D8:59:ALA:HB2	42:D8:96:ILE:HD13	1.80	0.64
35:68:2:ILE:HD12	35:68:6:THR:HG21	1.79	0.64
1:13:1412:C:H2'	1:13:1413:A:C8	2.33	0.64
23:2K:20:G:C2	23:2K:58:A:N3	2.66	0.64
46:D5:53:ILE:HG22	46:D5:71:VAL:HG13	1.80	0.64
1:1G:1347:G:O2'	1:1G:1373:G:O6	2.13	0.64
12:3I:117:ARG:HB3	12:3I:122:THR:HB	1.80	0.64
1:1G:1133:G:N2	1:1G:1141:C:O2	2.31	0.64
2:1E:98:LEU:HB2	2:1E:101:MET:HG3	1.79	0.64
1:13:1122:U:O4	1:13:1123:A:N6	2.31	0.64
40:B8:99:LEU:HB3	40:B8:101:PHE:CE1	2.33	0.64
12:3A:89:ARG:HG3	12:3A:97:ARG:HG2	1.79	0.64
16:7A:26:ARG:HH21	16:7A:31:LYS:HD3	1.62	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1G:1316:G:H4'	14:5A:18:VAL:HG11	1.79	0.64
1:13:345:C:O2'	1:13:346:G:N2	2.31	0.64
26:1H:2811:G:H1	26:1H:2889:C:H42	1.46	0.64
1:1G:1251:A:H2'	1:1G:1252:A:C8	2.33	0.64
24:3K:19:G:O2'	24:3K:57:G:N3	2.31	0.64
1:13:1301:U:O2'	1:13:1302:U:H5'	1.98	0.64
40:75:23:ARG:HG3	40:75:120:ARG:NH1	2.13	0.64
45:C5:8:LYS:HG3	45:C5:95:LYS:HE3	1.80	0.64
26:1H:543:C:H42	26:1H:550:G:H1	1.44	0.64
1:13:601:C:H2'	1:13:602:A:H8	1.62	0.64
1:1G:978:A:O2'	1:1G:1322:C:N3	2.31	0.64
44:F8:55:ASN:HB2	44:F8:80:ILE:HG13	1.79	0.64
26:1H:2080:G:H8	26:1H:2080:G:H5''	1.63	0.64
26:14:1264:G:OP1	52:J5:19:ARG:NH2	2.25	0.64
26:1H:1614:A:H2	58:1H:3859:HOH:O	1.80	0.63
49:G5:53:LEU:O	49:G5:57:ILE:HG13	1.97	0.63
2:12:43:ASP:O	2:12:47:THR:OG1	2.16	0.63
26:1H:2773:C:H5''	29:21:164:ARG:HG2	1.80	0.63
28:19:148:GLU:HB2	28:19:151:LYS:HD2	1.80	0.63
53:O8:14:THR:HA	53:O8:21:TYR:CD2	2.33	0.63
2:1E:126:GLU:HA	2:1E:129:GLU:HG2	1.80	0.63
1:13:324:G:N2	1:13:327:A:OP2	2.31	0.63
33:61:115:ALA:O	33:61:117:GLU:N	2.31	0.63
30:31:96:ASP:OD1	30:31:98:SER:HB3	1.98	0.63
26:14:1036:G:OP1	32:59:59:ARG:N	2.28	0.63
1:13:736:C:H2'	1:13:737:A:H8	1.64	0.63
1:13:1160:G:H1	1:13:1177:G:N2	1.94	0.63
26:14:2297:C:H2'	26:14:2298:A:H8	1.63	0.63
29:21:128:SER:OG	29:21:129:HIS:N	2.30	0.63
26:14:1342:A:H2	26:14:1602:U:H3	1.46	0.63
2:1E:212:GLN:O	2:1E:216:SER:OG	2.11	0.63
26:1H:743:G:O6	58:1H:3889:HOH:O	2.11	0.63
13:4I:10:PRO:HB2	13:4I:18:ALA:HB1	1.79	0.63
26:1H:121:G:OP1	58:1H:3996:HOH:O	2.15	0.63
55:Q8:9:GLY:H	55:Q8:12:LYS:HG3	1.62	0.63
1:13:963:G:H21	10:1I:55:LYS:CE	2.11	0.63
26:14:2250:G:C6	37:45:82:ARG:HD2	2.33	0.63
1:1G:1288:A:H4'	21:1B:13:ILE:HD13	1.80	0.63
38:98:55:ALA:HA	38:98:80:PHE:HE1	1.64	0.63
26:1H:443:A:H1'	26:1H:1201:C:O4'	1.98	0.63
1:1G:345:C:O3'	40:75:41:ARG:NH2	2.31	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:1H:831:G:N2	36:78:53:GLY:O	2.30	0.63
26:14:1966:A:H4'	26:14:1967:C:OP1	1.96	0.63
26:1H:587:C:OP2	36:78:21:ARG:NH2	2.30	0.63
2:12:178:ARG:NH1	2:12:196:LEU:O	2.29	0.63
34:58:96:GLU:O	34:58:98:VAL:HG12	1.99	0.63
38:55:24:GLN:NE2	38:55:36:THR:HG21	2.14	0.63
1:1G:280:C:H3'	1:1G:281:G:H5'	1.79	0.63
31:49:124:SER:HB2	31:49:131:TYR:CE1	2.33	0.63
30:31:33:LEU:HD23	36:78:1:MET:HG3	1.79	0.63
26:1H:2150:U:H2'	26:1H:2151:G:C8	2.32	0.63
44:B5:49:VAL:HB	44:B5:83:VAL:HG21	1.81	0.63
26:14:363(C):G:H2'	26:14:363(D):G:H8	1.63	0.63
13:4A:97:PRO:HB2	13:4A:101:GLN:HG3	1.81	0.63
14:5I:26:ARG:NH1	14:5I:43:CYS:SG	2.72	0.63
30:31:6:VAL:HG11	30:31:119:ARG:HA	1.80	0.63
26:1H:2611:U:H2'	52:N8:3:LYS:HG3	1.79	0.63
52:N8:50:GLY:H	52:N8:56:LYS:HG3	1.64	0.63
1:13:1240:U:OP2	7:6E:116:ALA:N	2.31	0.63
1:1G:588:G:H1	1:1G:651:C:N4	1.90	0.63
26:1H:2334:G:H5'	39:A8:9:ARG:HG2	1.80	0.63
26:1H:2371:G:H4'	53:O8:45:LYS:HG3	1.81	0.63
47:I8:53:MET:HG3	47:I8:59:LEU:HD23	1.80	0.63
26:14:2238:G:N7	58:14:3617:HOH:O	2.30	0.63
15:6I:6:GLU:HA	15:6I:9:GLN:HB2	1.81	0.63
26:14:972:G:OP2	26:14:973:A:O2'	2.09	0.63
2:1E:174:VAL:HG13	2:1E:184:VAL:HG11	1.79	0.63
1:13:649:G:H2'	1:13:650:G:H8	1.64	0.63
1:13:1259:C:N4	1:13:1260:C:O2	2.31	0.63
33:69:77:LEU:HD13	33:69:141:LYS:HB3	1.81	0.63
1:1G:1502:A:H5''	1:1G:1504:G:N7	2.13	0.63
26:14:567:A:P	58:14:3841:HOH:O	2.54	0.63
3:22:6:HIS:HB3	14:5A:49:HIS:CD2	2.33	0.63
29:29:200:GLU:N	29:29:200:GLU:OE2	2.25	0.63
2:12:68:ILE:HG12	2:12:161:ALA:HB3	1.80	0.63
26:14:4:C:H42	26:14:2899:G:H22	1.44	0.63
26:14:1664:A:OP2	58:14:3595:HOH:O	2.16	0.63
3:22:134:ILE:HG13	3:22:153:VAL:HG21	1.78	0.63
23:2L:10:G:N2	23:2L:27:G:H1'	2.14	0.63
26:1H:125:G:H5'	26:1H:125:G:H8	1.63	0.63
1:1G:1343:G:H2'	1:1G:1344:C:C6	2.33	0.63
1:1G:434:U:H2'	1:1G:435:C:C6	2.33	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:13:642:A:N3	8:7E:113:SER:OG	2.31	0.63
26:1H:1430:C:H2'	26:1H:1431:U:C6	2.32	0.63
26:1H:2572:A:N7	29:21:145:LYS:HB2	2.13	0.63
26:14:996:A:N6	26:14:1160:G:C6	2.67	0.63
1:13:1177:G:O2'	1:13:1178:G:O4'	2.16	0.63
26:14:2128:C:H42	26:14:2160:G:H1	1.45	0.63
26:1H:1159:U:OP1	50:L8:30:ARG:NH2	2.31	0.63
46:D5:4:ARG:NH1	46:D5:60:GLU:OE2	2.32	0.63
26:1H:2287:A:H62	26:1H:2344:U:H3	1.46	0.63
30:31:6:VAL:N	30:31:24:LEU:O	2.31	0.63
26:1H:2343:C:O2'	26:1H:2373:G:O2'	2.16	0.63
4:3E:173:TRP:CD1	4:3E:174:LEU:HG	2.33	0.63
4:32:173:TRP:CZ3	4:32:193:ASP:HB3	2.34	0.63
33:61:21:VAL:HG21	33:61:25:TYR:HD2	1.64	0.63
26:14:1778:U:H2'	26:14:1784:A:N6	2.14	0.63
45:C5:73:ARG:NH2	45:C5:81:LYS:O	2.30	0.63
2:12:22:LYS:HG2	2:12:40:HIS:NE2	2.14	0.63
5:4E:71:LEU:HD11	5:4E:114:GLY:HA3	1.81	0.63
29:21:167:VAL:HG21	29:21:187:ALA:HB3	1.80	0.63
26:1H:2311:A:H8	31:41:88:ILE:HD12	1.63	0.63
24:3K:75:C:HO2'	24:3K:76:A:H2	1.47	0.63
42:95:35:LEU:HB2	42:95:37:VAL:HG13	1.80	0.63
44:F8:3:THR:OG1	44:F8:4:ALA:HA	1.99	0.63
27:1J:44:G:O2'	27:1J:47:C:N4	2.30	0.63
8:7E:86:ILE:HG22	8:7E:87:SER:H	1.63	0.63
32:59:119:GLU:O	32:59:140:LYS:NZ	2.20	0.63
1:13:79:G:H21	1:13:89:U:H2'	1.63	0.63
42:95:48:GLY:H	42:95:52:VAL:HG22	1.64	0.63
1:1G:426:G:H4'	4:32:42:GLN:HA	1.81	0.63
37:45:98:LYS:HB3	37:45:99:PRO:HD2	1.81	0.63
1:1G:114:U:H2'	1:1G:115:G:C8	2.34	0.63
2:12:74:LYS:NZ	2:12:166:ASP:OD2	2.24	0.63
1:1G:420:U:O2'	1:1G:423:G:O6	2.09	0.63
50:L8:10:LYS:NZ	50:L8:15:TYR:OH	2.29	0.63
10:1A:8:LEU:HD22	10:1A:20:ALA:HB2	1.81	0.63
26:14:1364:G:OP2	48:F5:2:SER:N	2.32	0.63
24:1L:54:U:H5'	37:45:51:ARG:HH22	1.64	0.63
26:14:1688:U:O2	26:14:1700:A:H5'	1.99	0.63
55:Q8:46:ARG:HH21	55:Q8:48:PHE:HA	1.64	0.62
26:1H:563:G:OP2	58:1H:3639:HOH:O	2.16	0.62
26:1H:860:U:C5	26:1H:917:A:H2	2.15	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:32:26:CYS:HA	4:32:31:CYS:HA	1.81	0.62
31:49:66:GLN:OE1	31:49:98:ARG:NH1	2.31	0.62
26:1H:1796:U:H2'	26:1H:1797:C:C6	2.33	0.62
26:14:1784:A:H5''	58:14:3559:HOH:O	1.97	0.62
37:45:35:VAL:HG22	37:45:102:VAL:HG22	1.82	0.62
26:14:2015:A:N3	52:J5:2:ALA:N	2.46	0.62
26:1H:155:C:N4	26:1H:171:G:H1	1.87	0.62
1:1G:957:U:H2'	1:1G:959:A:OP2	1.99	0.62
12:3I:93:LEU:O	12:3I:96:VAL:HG13	1.99	0.62
9:82:112:LYS:HG2	9:82:119:ALA:HB2	1.81	0.62
4:3E:31:CYS:SG	4:3E:32:ALA:N	2.72	0.62
13:4A:57:ARG:NH1	51:I5:34:GLU:O	2.31	0.62
26:14:824:A:H1'	26:14:2358:G:N7	2.14	0.62
13:4I:108:ARG:NH1	13:4I:112:GLY:O	2.31	0.62
1:13:667:G:H4'	15:6I:51:HIS:CE1	2.34	0.62
26:1H:270(L):U:H3	33:61:50:ARG:HG2	1.64	0.62
42:D8:1:MET:HG2	42:D8:43:GLU:HB3	1.80	0.62
11:2A:48:ILE:HD11	11:2A:64:ALA:HA	1.81	0.62
6:52:2:ARG:HD3	6:52:92:LYS:HE3	1.81	0.62
36:78:19:VAL:HB	36:78:27:HIS:HB2	1.79	0.62
24:3K:6:G:N2	24:3K:68:C:N3	2.47	0.62
1:13:600:C:H2'	1:13:601:C:C6	2.34	0.62
1:13:601:C:H2'	1:13:602:A:C8	2.33	0.62
26:14:1013:C:H42	26:14:1149:G:H1	1.45	0.62
9:8E:26:VAL:HG13	9:8E:61:ALA:HB3	1.80	0.62
16:7A:34:GLU:OE2	16:7A:55:ARG:NH1	2.32	0.62
1:13:1429:C:H2'	1:13:1430:C:H6	1.65	0.62
13:4A:34:LEU:O	13:4A:38:GLY:N	2.32	0.62
1:1G:711:G:OP1	6:52:54:LYS:NZ	2.29	0.62
44:B5:36:LYS:HG3	44:B5:56:THR:HG23	1.81	0.62
26:1H:2600:A:N6	58:1H:3678:HOH:O	2.22	0.62
26:14:2012:G:H5''	43:A5:96:ILE:HD11	1.81	0.62
49:K8:15:LYS:H	49:K8:67:LYS:HZ3	1.46	0.62
2:1E:179:LYS:HA	8:7E:72:PRO:HG3	1.82	0.62
38:98:55:ALA:HA	38:98:80:PHE:CE1	2.33	0.62
1:1G:1141:C:H2'	1:1G:1142:G:H8	1.64	0.62
26:14:801:G:OP2	58:14:3700:HOH:O	2.15	0.62
7:62:92:SER:HB2	7:62:94:ARG:HG2	1.82	0.62
26:14:2378:A:H4'	39:65:23:ARG:HH11	1.64	0.62
32:59:42:ARG:NH1	32:59:53:GLU:O	2.33	0.62
4:3E:102:ASP:OD1	4:3E:103:ASN:N	2.32	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1G:1054:C:O2'	1:1G:1055:A:O5'	2.14	0.62
33:69:80:PRO:HA	33:69:143:SER:HA	1.79	0.62
1:1G:1502:A:H4'	1:1G:1503:A:OP2	1.99	0.62
26:14:259:G:H21	26:14:621:A:H8	1.48	0.62
1:1G:1376:U:H2'	1:1G:1377:A:C8	2.34	0.62
26:1H:736:C:H5''	58:1H:4337:HOH:O	1.98	0.62
39:A8:51:ALA:HB3	39:A8:73:LEU:HG	1.81	0.62
10:1I:4:ILE:HG12	10:1I:100:THR:HA	1.82	0.62
24:3L:72:C:H3'	24:3L:73:A:H5''	1.81	0.62
26:14:329:G:P	45:C5:71:LYS:HD3	2.39	0.62
1:1G:1047:G:H1	1:1G:1210:C:H42	1.47	0.62
48:J8:87:PRO:HA	48:J8:90:ILE:HG13	1.81	0.62
1:1G:1060:C:H5''	10:1A:51:ARG:HG2	1.80	0.62
34:58:96:GLU:HG2	34:58:97:ARG:N	2.13	0.62
1:13:390:C:O3'	16:7I:28:ARG:NH2	2.31	0.62
26:1H:1858:G:H2'	26:1H:1883:G:H22	1.64	0.62
11:2A:32:ILE:HD11	11:2A:68:ALA:HB1	1.81	0.62
1:1G:222:U:H2'	1:1G:223:U:C6	2.35	0.62
7:62:84:ASN:OD1	7:62:84:ASN:N	2.32	0.62
26:14:1729:A:H2'	26:14:1731:G:N2	2.15	0.62
26:1H:1514:U:H2'	26:1H:1515:C:C6	2.33	0.62
45:C5:18:GLY:O	45:C5:20:TYR:N	2.32	0.62
1:1G:1002:G:H2'	1:1G:1003:G:H8	1.63	0.62
29:29:89:ASP:OD1	29:29:90:THR:N	2.33	0.62
12:3A:78:GLN:HG3	12:3A:81:SER:HB2	1.80	0.62
26:14:1485:G:N2	26:14:1504:C:N3	2.43	0.62
26:1H:1534:G:H2'	26:1H:1535:U:H4'	1.81	0.62
30:39:7:TYR:HE2	30:39:10:PRO:HG3	1.65	0.62
33:69:75:LEU:HD22	33:69:76:THR:H	1.65	0.62
30:39:63:LYS:HE2	30:39:67:GLN:HB3	1.80	0.62
47:E5:26:TYR:O	47:E5:29:GLN:HB2	2.00	0.62
32:59:101:ARG:NH1	32:59:121:ILE:O	2.33	0.62
1:13:1061:G:OP1	10:1I:59:SER:OG	2.15	0.62
1:1G:980:C:H3'	1:1G:981:U:C6	2.34	0.62
1:1G:716:A:N3	11:2A:118:GLY:HA2	2.15	0.62
6:52:97:PHE:HD2	18:9A:31:LEU:HD21	1.65	0.62
26:1H:1510:A:O2'	26:1H:1512:G:N7	2.27	0.62
22:1K:76:A:O3'	26:1H:2506:U:H1'	1.98	0.62
32:59:6:ARG:HH11	32:59:6:ARG:N	1.98	0.62
26:14:67:U:H2'	26:14:68:G:C8	2.34	0.62
13:4A:8:GLU:OE1	13:4A:9:ILE:N	2.32	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
37:45:135:ASP:OD2	46:D5:81:ARG:NH2	2.33	0.62
1:13:7:G:H5'	1:13:298:A:O4'	1.99	0.62
1:13:624:C:O3'	16:7I:10:GLY:HA2	1.99	0.62
15:6A:17:ARG:HD3	15:6A:26:GLU:HG3	1.82	0.62
26:1H:2584:U:H2'	26:1H:2585:U:H2'	1.82	0.62
40:75:105:LEU:HD23	40:75:109:GLU:HG3	1.80	0.62
17:8I:31:LEU:HD23	17:8I:32:TYR:CZ	2.35	0.62
26:14:2153:G:N2	26:14:2154:G:O6	2.32	0.62
7:62:27:ILE:HA	7:62:30:ILE:HD12	1.82	0.62
1:1G:1300:G:O2'	1:1G:1301:U:O5'	2.13	0.62
2:1E:11:LEU:O	2:1E:16:HIS:NE2	2.33	0.62
47:E5:18:ALA:HB3	47:E5:20:ARG:HH21	1.64	0.62
36:78:50:ARG:HD3	55:Q8:7:HIS:CD2	2.35	0.62
9:82:112:LYS:HA	9:82:119:ALA:HB2	1.81	0.62
26:1H:1178:C:H4'	26:1H:1179:C:OP1	2.00	0.62
26:1H:1711:C:H2'	26:1H:1712:C:H6	1.64	0.62
26:1H:2314:C:H2'	26:1H:2315:G:H8	1.65	0.62
19:AI:63:THR:OG1	19:AI:64:GLU:N	2.32	0.62
3:2E:11:ARG:HH21	3:2E:180:ALA:HB3	1.64	0.62
26:14:2150:U:H2'	26:14:2151:G:H8	1.64	0.62
28:11:108:PRO:HG3	28:11:143:HIS:CE1	2.34	0.62
26:1H:1188:U:H4'	42:D8:79:VAL:HG22	1.82	0.61
26:1H:2503:A:H4'	26:1H:2504:U:OP1	2.00	0.61
24:3K:4:C:H2'	24:3K:5:G:C8	2.35	0.61
32:51:6:ARG:HA	32:51:66:GLY:HA2	1.82	0.61
13:4A:91:ARG:NH2	13:4A:97:PRO:O	2.33	0.61
26:14:93:C:H5'	26:14:94:G:OP2	2.00	0.61
1:1G:1274:G:H2'	1:1G:1275:A:C8	2.35	0.61
26:14:1678:G:H22	26:14:1989:G:N2	1.98	0.61
26:14:303:U:H2'	26:14:304:G:C8	2.35	0.61
26:1H:848:G:H2'	26:1H:849:A:C8	2.35	0.61
46:H8:151:HIS:HB3	46:H8:168:GLU:HA	1.81	0.61
26:1H:1509:C:H3'	26:1H:1510:A:H5''	1.82	0.61
26:1H:125:G:H5'	26:1H:125:G:C8	2.35	0.61
37:45:20:ALA:HA	37:45:99:PRO:HG2	1.83	0.61
1:1G:971:G:N2	1:1G:1363:A:OP2	2.33	0.61
39:A8:52:SER:HB2	39:A8:55:ALA:H	1.65	0.61
26:14:528:A:O2'	26:14:529:A:H5'	2.00	0.61
1:1G:1423:G:H2'	1:1G:1424:C:C6	2.35	0.61
26:1H:581:C:H2'	26:1H:582:G:C8	2.35	0.61
28:19:44:ASN:ND2	28:19:46:GLN:HG3	2.15	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:13:811:C:N3	58:13:1805:HOH:O	2.31	0.61
26:1H:323:G:C8	30:31:171:PRO:HG3	2.35	0.61
26:14:1819:A:H4'	26:14:1820:U:O5'	2.00	0.61
26:1H:2505:G:O6	26:1H:2576:G:H2'	2.00	0.61
1:1G:1279:A:O2'	1:1G:1282:C:N4	2.33	0.61
15:6I:17:ARG:HH11	15:6I:17:ARG:HG3	1.65	0.61
26:1H:1316:U:H2'	26:1H:1317:A:C8	2.33	0.61
4:32:18:LYS:HB3	4:32:33:MET:HG3	1.81	0.61
26:14:2689:U:H4'	26:14:2690:C:H5'	1.82	0.61
1:1G:745:C:H5'	1:1G:851:G:H21	1.63	0.61
37:45:21:THR:HG22	37:45:23:GLY:HA3	1.81	0.61
26:1H:779:U:OP1	28:11:49:ILE:HG13	2.00	0.61
1:13:453:A:H4'	16:7I:72:ARG:HB2	1.81	0.61
26:1H:654(J):A:N1	26:1H:654(M):C:N4	2.48	0.61
41:C8:11:ARG:O	41:C8:15:LYS:HG3	2.01	0.61
26:14:2210:G:H3'	26:14:2211:G:N2	2.16	0.61
16:7I:50:LYS:HD3	16:7I:51:VAL:H	1.63	0.61
26:1H:2250:G:C4	37:88:82:ARG:HG2	2.35	0.61
1:13:1028:C:H42	1:13:1033:G:H1	1.48	0.61
3:22:8:ILE:HG23	3:22:16:ARG:HG2	1.83	0.61
26:14:2675:A:H4'	35:25:29:ASN:HD21	1.65	0.61
26:1H:2347:C:O5'	53:O8:39:TYR:OH	2.11	0.61
32:51:157:TYR:O	32:51:158:HIS:ND1	2.32	0.61
1:13:1060:C:HO2'	10:1I:56:HIS:HD1	1.41	0.61
38:98:9:LYS:HA	38:98:17:ARG:NE	2.15	0.61
32:51:113:VAL:HG11	32:51:151:ILE:HD13	1.83	0.61
49:K8:32:LEU:HD11	49:K8:54:LYS:HG3	1.82	0.61
29:21:54:GLN:O	29:21:55:ASN:ND2	2.33	0.61
42:95:85:LYS:HG3	42:95:87:HIS:H	1.64	0.61
1:13:1182:G:H4'	1:13:1183:A:H5''	1.81	0.61
4:32:9:CYS:SG	4:32:22:LYS:HD2	2.41	0.61
37:88:51:ARG:HH11	37:88:51:ARG:HB3	1.64	0.61
29:21:119:ARG:HG3	29:21:119:ARG:HH11	1.65	0.61
26:14:2572:A:OP1	26:14:2574:G:O2'	2.18	0.61
26:14:2016:U:O2	52:J5:7:PRO:HG2	2.00	0.61
1:1G:1286:A:H3'	1:1G:1286:A:H8	1.65	0.61
1:1G:1423:G:H2'	1:1G:1424:C:H6	1.65	0.61
49:K8:42:GLY:O	49:K8:44:LEU:N	2.33	0.61
37:88:12:GLN:HG2	37:88:73:PRO:HD2	1.83	0.61
11:2A:27:ASN:OD1	11:2A:28:THR:N	2.33	0.61
26:14:1412:A:H2'	26:14:1413:G:C8	2.35	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:8E:89:ASN:O	9:8E:89:ASN:ND2	2.30	0.61
26:1H:952:G:H5''	26:1H:953:A:OP2	2.00	0.61
26:14:322:A:H3'	30:39:169:ASN:OD1	2.00	0.61
26:1H:482:A:OP2	26:1H:482:A:H8	1.83	0.61
48:F5:84:GLY:O	48:F5:87:PRO:HD2	2.00	0.61
1:1G:1349:A:P	9:82:118:LYS:HZ1	2.22	0.61
1:13:1452:C:O2'	1:13:1453:G:OP2	2.18	0.61
11:2A:82:VAL:HG13	11:2A:108:ILE:HG23	1.83	0.61
26:14:7:G:H2'	26:14:8:A:C8	2.35	0.61
1:1G:567:G:N3	58:1G:1738:HOH:O	2.31	0.61
26:1H:2849:U:H4'	26:1H:2868:A:C2	2.35	0.61
34:15:56:ASN:HA	34:15:125:GLY:H	1.64	0.61
20:BI:49:ALA:CB	20:BI:99:LEU:HB2	2.31	0.61
26:14:994:C:O2'	26:14:996:A:OP1	2.19	0.61
1:13:1178:G:H5''	9:8E:93:ARG:HH22	1.66	0.61
26:14:1856:G:H1	26:14:1886:C:N4	1.99	0.61
12:3A:41:ARG:HH12	12:3A:43:VAL:HG22	1.65	0.61
1:1G:503:C:OP2	12:3A:116:SER:HB3	2.00	0.61
26:14:588:U:H2'	26:14:589:C:C6	2.36	0.61
1:1G:517:G:N2	1:1G:530:G:OP1	2.21	0.61
26:14:990:A:H8	26:14:990:A:H5'	1.65	0.61
36:78:52:GLU:HG3	36:78:57:THR:HA	1.83	0.61
41:C8:97:ASP:OD2	41:C8:101:ARG:NH1	2.34	0.61
36:78:18:ARG:O	36:78:19:VAL:HG22	2.00	0.61
27:1J:16:G:H2'	27:1J:17:C:C6	2.36	0.61
26:1H:1509:C:H2'	26:1H:1511:A:C8	2.34	0.61
24:3K:19:G:C6	26:1H:2112:G:H1'	2.36	0.61
26:1H:1814:G:C6	58:1H:4231:HOH:O	2.51	0.61
13:4A:86:CYS:SG	13:4A:88:ARG:HB3	2.40	0.61
1:13:757:U:H2'	1:13:758:G:O4'	2.00	0.61
26:1H:639:U:H2'	26:1H:640:C:C6	2.35	0.61
1:13:881:G:OP2	12:3I:12:ARG:NH2	2.33	0.61
26:1H:2887:U:H2'	26:1H:2888:C:C6	2.35	0.61
26:14:446:G:OP2	58:14:3676:HOH:O	2.16	0.61
3:22:36:ASP:HA	3:22:39:ILE:HD12	1.82	0.61
46:D5:11:GLU:CD	46:D5:12:GLY:H	2.03	0.61
1:1G:1126:U:N3	1:1G:1281:U:O4'	2.34	0.61
26:14:1022:G:H22	26:14:1142(A):A:H2	1.48	0.61
47:E5:49:LYS:HE3	47:E5:82:ARG:NH1	2.16	0.61
26:14:900:A:H3'	26:14:901:A:H8	1.65	0.61
3:22:88:ARG:HA	3:22:101:LEU:HD13	1.81	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
40:75:26:ASP:O	40:75:49:VAL:HG22	2.01	0.61
1:13:484:G:O2'	1:13:485:G:OP2	2.17	0.61
26:14:1664:A:OP2	58:14:3594:HOH:O	2.16	0.61
26:14:1678:G:H22	26:14:1989:G:H22	1.48	0.61
26:14:1198:U:H2'	26:14:1199:U:C6	2.36	0.61
3:22:70:VAL:HG12	3:22:72:LYS:H	1.65	0.61
1:13:591:U:H2'	1:13:592:G:H8	1.65	0.61
39:A8:27:SER:HA	39:A8:88:ASP:HB3	1.82	0.61
24:3L:33:U:O2'	24:3L:35:A:OP2	2.18	0.61
26:1H:2702:U:H6	26:1H:2702:U:OP1	1.84	0.61
1:1G:926:G:N2	25:4L:15:A:OP2	2.33	0.61
26:1H:49:A:N7	26:1H:120:U:C5	2.64	0.61
40:B8:26:ASP:O	40:B8:49:VAL:HG12	2.00	0.61
26:1H:1287:A:C8	38:98:107:ASP:HB2	2.36	0.61
43:E8:17:VAL:HG13	43:E8:76:VAL:HG11	1.82	0.61
3:2E:44:GLU:HA	3:2E:52:LEU:HD11	1.82	0.61
26:14:2273:A:H2'	26:14:2274:A:H8	1.66	0.61
26:14:2845:G:H5''	40:75:54:ARG:O	2.00	0.61
10:1A:8:LEU:HA	10:1A:96:ILE:HG22	1.82	0.61
26:1H:910:A:N7	37:88:13:GLN:HG3	2.15	0.61
1:1G:371:G:O2'	1:1G:373:A:N7	2.33	0.61
26:1H:66:C:H2'	26:1H:67:U:H6	1.66	0.61
26:1H:919:G:N2	26:1H:2269:A:OP2	2.33	0.61
16:7A:75:ARG:HG3	16:7A:80:PHE:CD2	2.36	0.61
44:B5:51:VAL:H	44:B5:83:VAL:HG23	1.65	0.60
1:13:110:C:H2'	1:13:111:G:O4'	2.01	0.60
26:14:2611:U:H3'	26:14:2611:U:OP2	2.01	0.60
26:14:736:C:OP1	58:14:3934:HOH:O	2.17	0.60
30:31:129:PHE:HA	30:31:142:TRP:NE1	2.16	0.60
35:25:102:VAL:HB	35:25:106:LEU:HD12	1.81	0.60
51:I5:18:CYS:H	51:I5:19:GLY:HA2	1.66	0.60
11:2A:100:ALA:O	11:2A:102:GLY:N	2.34	0.60
12:3I:59:ARG:HA	12:3I:65:GLU:HA	1.83	0.60
1:13:1305:G:H21	1:13:1331:G:H2'	1.66	0.60
53:K5:10:LEU:HB3	55:M5:34:TRP:CZ3	2.36	0.60
26:14:654(E):C:N4	26:14:654(P):G:H22	1.99	0.60
24:3L:3:C:H2'	24:3L:4:C:O4'	2.01	0.60
1:13:1060:C:H5''	10:1I:51:ARG:HG2	1.83	0.60
26:1H:2849:U:OP2	40:B8:95:ARG:NH1	2.34	0.60
19:AI:19:VAL:HG13	19:AI:47:HIS:HD2	1.65	0.60
26:1H:547:A:H2'	26:1H:548:A:C8	2.35	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:13:1497:G:H2'	1:13:1498:U:H5'	1.82	0.60
26:1H:2243:U:OP1	58:1H:4129:HOH:O	2.16	0.60
11:2A:18:ARG:NH2	11:2A:35:PRO:O	2.34	0.60
26:14:155:C:H42	26:14:171:G:H1	1.49	0.60
32:51:8:PRO:HG2	32:51:69:ARG:NH2	2.15	0.60
11:2A:67:ASP:OD2	11:2A:71:LYS:NZ	2.34	0.60
24:3K:58:A:O2'	24:3K:59:U:OP1	2.19	0.60
26:14:2275:C:H6	26:14:2275:C:H5'	1.67	0.60
27:1J:56:G:H4'	27:1J:57:A:C8	2.37	0.60
1:1G:728:A:H2'	1:1G:729:A:C8	2.35	0.60
45:G8:55:TYR:HB2	45:G8:58:GLY:HA3	1.81	0.60
26:1H:880:G:O2'	26:1H:881:G:O5'	2.19	0.60
1:1G:1028(A):C:O2	1:1G:1033:G:N2	2.34	0.60
1:1G:1095:U:OP1	1:1G:1108:G:N1	2.35	0.60
30:31:108:LYS:O	30:31:112:MET:HG3	2.00	0.60
33:69:54:GLN:HA	33:69:57:ARG:HB3	1.84	0.60
16:7A:37:GLY:HA2	16:7A:50:LYS:HD3	1.82	0.60
26:1H:957:A:N1	26:1H:2458:G:H4'	2.16	0.60
26:1H:2817:G:H1	26:1H:2829:C:H42	1.50	0.60
1:13:1007:C:N4	1:13:1022:G:H1	1.99	0.60
26:1H:140:A:H8	26:1H:1408:C:HO2'	1.48	0.60
26:14:2280:G:O2'	26:14:2388:A:N1	2.33	0.60
26:1H:1668:A:N6	26:1H:1676:A:H61	1.99	0.60
51:I5:20:ASN:OD1	51:I5:36:CYS:HB2	2.02	0.60
40:B8:16:ARG:HE	40:B8:19:LEU:HD11	1.65	0.60
9:82:10:ARG:HD2	9:82:105:ASP:HB3	1.82	0.60
26:1H:1086:A:H1'	26:1H:1103:A:H61	1.64	0.60
26:1H:176:G:O2'	26:1H:177:G:H5'	2.02	0.60
26:1H:2857:G:N2	26:1H:2860:A:OP2	2.27	0.60
31:49:173:LEU:HA	31:49:176:LEU:HB2	1.83	0.60
3:22:14:ILE:HG12	3:22:15:THR:H	1.66	0.60
36:78:50:ARG:HD3	55:Q8:7:HIS:NE2	2.16	0.60
45:G8:76:CYS:HB2	45:G8:82:PRO:HD3	1.83	0.60
28:11:223:GLY:HA3	28:11:231:HIS:CE1	2.36	0.60
26:1H:2165:G:N7	26:1H:2166:G:N2	2.48	0.60
37:45:75:THR:HA	37:45:89:ASN:HA	1.84	0.60
26:1H:873:G:N2	26:1H:904:C:O2	2.27	0.60
30:31:114:VAL:HG11	30:31:202:PHE:HE2	1.67	0.60
15:6A:43:LEU:HD11	15:6A:53:HIS:HA	1.84	0.60
46:H8:7:ALA:HB3	46:H8:61:LEU:HB2	1.84	0.60
26:14:1794:U:O2'	26:14:1795:C:H5'	2.00	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
13:4I:27:LYS:HA	13:4I:31:LYS:HZ3	1.65	0.60
26:1H:1230:C:H2'	26:1H:1231:G:C8	2.37	0.60
26:1H:1607:C:H4'	26:1H:1608:A:O5'	2.01	0.60
8:72:7:ALA:O	8:72:11:THR:OG1	2.12	0.60
26:1H:620:G:H4'	26:1H:621:A:C5'	2.31	0.60
26:14:1757:U:N3	26:14:1762:A:H2	1.96	0.60
26:1H:399:G:OP2	58:1H:4238:HOH:O	2.16	0.60
36:78:58:THR:HG21	55:Q8:52:LYS:HD3	1.83	0.60
26:1H:1534:G:H22	26:1H:1538:G:H22	1.48	0.60
49:K8:58:ALA:O	49:K8:62:THR:HG22	2.00	0.60
35:25:59:LYS:HB3	35:25:87:ILE:HG22	1.83	0.60
28:11:68:LYS:HB3	28:11:70:TRP:CH2	2.36	0.60
5:4E:77:PRO:HD2	5:4E:142:LEU:HD13	1.83	0.60
26:14:2836:U:H2'	26:14:2837:G:C8	2.37	0.60
1:1G:976:G:OP2	1:1G:1358:U:O2'	2.18	0.60
26:1H:2598:A:OP1	58:1H:3655:HOH:O	2.16	0.60
3:22:11:ARG:NH1	3:22:11:ARG:HB2	2.17	0.60
8:7E:41:ARG:NH2	8:7E:123:GLU:OE1	2.33	0.60
1:13:1306:A:H61	1:13:1331:G:H1'	1.66	0.60
26:1H:2347:C:P	53:O8:39:TYR:HH	2.24	0.60
53:K5:18:ARG:NH2	53:K5:43:CYS:O	2.34	0.60
26:14:2335:A:O2'	26:14:2337:G:N7	2.31	0.60
26:1H:534:U:H5'	41:C8:42:ALA:HB1	1.83	0.60
26:14:870:A:H2'	26:14:871:U:O4'	2.02	0.60
10:1A:40:LEU:HD12	10:1A:41:PRO:HD2	1.83	0.60
34:58:38:HIS:O	41:C8:67:ALA:HB1	2.01	0.60
46:D5:139:VAL:HG13	46:D5:156:LYS:HE2	1.82	0.60
39:65:30:ARG:HG3	39:65:35:ILE:HD13	1.83	0.60
31:49:32:PRO:HB2	31:49:172:LEU:HD13	1.83	0.60
18:9A:21:LYS:HZ1	18:9A:57:GLY:HA3	1.66	0.60
46:H8:163:LEU:HD13	46:H8:165:VAL:HA	1.82	0.60
20:BI:26:ASN:HB2	20:BI:71:THR:HG23	1.83	0.60
1:1G:1298:C:H41	7:62:114:ARG:HA	1.67	0.60
1:13:1028(A):C:H42	1:13:1032(A):G:H1	1.50	0.60
32:51:54:ARG:HD3	32:51:65:HIS:ND1	2.16	0.60
1:1G:1240:U:N3	7:62:32:ARG:HG3	2.16	0.60
8:72:20:TYR:HA	8:72:65:TYR:CZ	2.37	0.60
24:3L:9:A:O2'	24:3L:10:G:N7	2.32	0.60
31:49:56:ALA:HA	31:49:59:GLU:HB2	1.84	0.60
30:31:160:ASN:OD1	30:31:163:VAL:HG23	2.02	0.60
1:1G:1099:G:OP1	2:12:148:TYR:OH	2.19	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
31:41:65:GLY:HA2	51:M8:7:PRO:HG2	1.84	0.60
1:1G:690:G:H2'	1:1G:691:G:O4'	2.01	0.60
1:1G:785:G:N7	58:1G:1746:HOH:O	2.31	0.60
26:14:729:G:P	28:19:208:LYS:HZ1	2.25	0.60
1:1G:1160:G:N7	1:1G:1181:G:N1	2.40	0.60
26:1H:1705:G:C2'	26:1H:1706:U:H5'	2.31	0.60
1:13:768:A:OP2	58:13:1885:HOH:O	2.17	0.60
12:3A:27:LEU:HG	12:3A:33:ARG:HG2	1.84	0.60
1:13:631:G:H2'	1:13:632:A:N3	2.16	0.60
12:3I:89:ARG:HD3	12:3I:91:LYS:HB3	1.83	0.60
1:13:276:G:O3'	17:8I:68:ARG:NH1	2.34	0.60
1:13:1263:C:H2'	1:13:1264:C:C6	2.36	0.60
12:3A:58:VAL:O	12:3A:65:GLU:HA	2.02	0.60
21:1B:8:THR:HG22	21:1B:11:GLY:H	1.65	0.60
26:1H:10:G:N2	26:1H:2801:A:O2'	2.35	0.60
1:13:236:G:H5''	17:8I:42:TYR:OH	2.02	0.60
26:1H:764:A:O4'	28:11:213:ARG:HG3	2.02	0.60
26:1H:607:U:N3	26:1H:621:A:C2	2.67	0.59
15:6I:16:ALA:HB1	15:6I:21:ASP:HB3	1.84	0.59
29:21:101:ARG:HG2	29:21:169:ASN:OD1	2.02	0.59
27:16:40:U:H5	51:M8:2:LYS:HE2	1.67	0.59
26:14:2261:C:O2'	26:14:2262:U:H5'	2.02	0.59
26:1H:1113:U:H2'	26:1H:1114:G:C8	2.36	0.59
26:1H:357:A:H2'	26:1H:358:U:C6	2.37	0.59
35:25:10:VAL:HG22	35:25:17:ARG:O	2.02	0.59
32:51:12:PRO:HG2	32:51:13:LYS:HG2	1.84	0.59
26:1H:1693:U:H1'	28:11:14:ARG:NH2	2.16	0.59
26:14:2123:G:H2'	26:14:2124:G:H8	1.67	0.59
26:14:861:A:N3	27:1J:79:C:O2'	2.34	0.59
37:88:133:ARG:O	37:88:134:ARG:HB2	2.01	0.59
1:1G:147:G:H1	1:1G:175:C:H42	1.50	0.59
32:51:144:VAL:O	32:51:148:ILE:HG12	2.02	0.59
19:AA:50:ALA:CB	19:AA:57:HIS:HB3	2.32	0.59
12:3A:41:ARG:NH1	12:3A:41:ARG:HB3	2.17	0.59
11:2A:32:ILE:HD13	11:2A:72:ALA:HB2	1.84	0.59
23:2L:20:G:C2	23:2L:58:A:C2	2.90	0.59
2:12:63:MET:HG3	2:12:225:ALA:HB1	1.84	0.59
26:14:2528:U:O2'	26:14:2530:A:OP1	2.16	0.59
37:88:21:THR:OG1	37:88:22:LYS:O	2.15	0.59
26:14:140:A:H8	26:14:1408:C:HO2'	1.46	0.59
3:22:21:ARG:HH12	10:1A:92:THR:HB	1.67	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:1366:A:H2'	26:14:1367:A:O4'	2.02	0.59
26:1H:58:G:N2	26:1H:70:G:C4	2.70	0.59
1:1G:556:C:OP2	12:3A:20:LYS:NZ	2.35	0.59
1:1G:1316:G:H5''	14:5A:17:LYS:NZ	2.17	0.59
26:14:2151:G:H2'	26:14:2152:G:H8	1.67	0.59
37:45:38:GLU:HB2	37:45:127:ILE:HG22	1.83	0.59
1:1G:632:A:H1'	1:1G:633:G:OP2	2.01	0.59
24:3L:15:G:H22	24:3L:48:C:H41	1.50	0.59
1:1G:45:U:H2'	1:1G:46:G:C8	2.37	0.59
26:1H:1374:G:H2'	26:1H:1375:C:H6	1.66	0.59
1:1G:581:G:OP1	15:6A:61:GLY:HA3	2.02	0.59
4:3E:74:GLN:O	4:3E:78:LEU:HD13	2.02	0.59
33:61:109:ILE:HB	33:61:130:TYR:CZ	2.36	0.59
26:14:172:C:H2'	26:14:173:G:H8	1.66	0.59
1:1G:1503:A:N3	25:4L:13:A:N6	2.51	0.59
26:14:1771:C:O2'	26:14:1786:A:H8	1.85	0.59
34:15:4:TYR:O	41:85:64:ARG:NH1	2.35	0.59
2:1E:178:ARG:NH1	2:1E:196:LEU:O	2.31	0.59
26:1H:1331:A:HO2'	26:1H:1332:G:H8	1.51	0.59
26:1H:2636:U:H3	26:1H:2782:G:H1	1.49	0.59
52:N8:41:PRO:HG2	52:N8:44:THR:HG21	1.82	0.59
18:9A:53:ARG:NE	18:9A:59:SER:O	2.23	0.59
31:41:143:GLU:OE1	51:M8:26:SER:OG	2.20	0.59
1:1G:683:G:C6	1:1G:684:A:C6	2.90	0.59
38:98:87:TYR:HD1	38:98:90:ARG:HD2	1.68	0.59
26:1H:1438:U:H2'	26:1H:1439:A:C8	2.37	0.59
28:11:59:LYS:HD2	28:11:60:ARG:N	2.16	0.59
26:1H:580:C:H2'	26:1H:581:C:C6	2.38	0.59
18:9A:56:THR:HB	18:9A:58:LEU:HD12	1.84	0.59
26:14:247:G:H4'	26:14:386:G:C5	2.37	0.59
1:13:142:G:H2'	1:13:143:A:H8	1.68	0.59
1:13:692:U:H5'	1:13:797:C:H5'	1.85	0.59
26:1H:625:G:N7	36:78:107:LYS:NZ	2.50	0.59
35:68:71:ARG:NH1	40:B8:74:ARG:HH21	2.00	0.59
26:14:1111:A:H5'	32:59:3:ARG:HD3	1.85	0.59
27:1J:24:G:H4'	27:1J:25:A:H5'	1.85	0.59
45:C5:39:VAL:O	45:C5:40:GLU:HB2	2.02	0.59
31:49:95:ARG:HG2	31:49:96:ARG:H	1.67	0.59
26:14:2712:U:H2'	26:14:2714:G:H5''	1.83	0.59
3:2E:40:ARG:HG3	3:2E:40:ARG:HH11	1.65	0.59
42:95:43:GLU:OE2	42:95:44:LYS:NZ	2.27	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:13:1429:C:H2'	1:13:1430:C:C6	2.36	0.59
26:14:288:C:O2	26:14:353:G:N2	2.27	0.59
40:75:56:GLY:O	40:75:59:THR:HG23	2.03	0.59
1:1G:1128:C:O2	1:1G:1144:G:N2	2.35	0.59
1:1G:1227:A:OP1	19:AA:80:TYR:OH	2.19	0.59
26:14:1266:G:O4'	43:A5:15:ARG:NH2	2.35	0.59
39:A8:34:HIS:NE2	39:A8:54:LEU:HD23	2.17	0.59
35:25:13:ASN:ND2	35:25:96:THR:HG23	2.17	0.59
1:1G:1203:C:H2'	1:1G:1204:A:O4'	2.03	0.59
3:2E:8:ILE:HD12	3:2E:16:ARG:HG3	1.83	0.59
26:1H:606:U:H4'	26:1H:658:C:H4'	1.85	0.59
1:13:272:C:H2'	1:13:273:A:C8	2.37	0.59
1:13:342:C:H2'	1:13:343:U:O4'	2.01	0.59
30:31:6:VAL:HG21	30:31:119:ARG:HB2	1.85	0.59
26:14:527:C:H4'	26:14:528:A:O5'	2.02	0.59
1:1G:1326:C:OP1	21:1B:17:THR:OG1	2.15	0.59
1:13:445:G:H1	1:13:489:C:H42	1.48	0.59
15:6I:42:HIS:HD2	15:6I:43:LEU:HD23	1.67	0.59
1:13:201:C:H42	1:13:216:G:H1	1.49	0.59
26:1H:654(A):A:H2	26:1H:654(T):A:N1	2.00	0.59
26:14:2781:A:H5''	26:14:2782:G:H5'	1.85	0.59
31:41:68:PRO:HB3	31:41:92:VAL:HB	1.85	0.59
38:98:1:MET:O	38:98:1:MET:HG2	2.02	0.59
14:5A:27:CYS:SG	14:5A:29:ARG:HG2	2.43	0.59
26:1H:1265:A:H3'	52:N8:19:ARG:HH12	1.68	0.59
1:1G:618:C:H5'	1:1G:619:U:H5''	1.84	0.59
50:L8:30:ARG:HH11	50:L8:33:GLN:HE21	1.51	0.59
1:1G:192:U:O4'	20:BA:103:GLY:HA2	2.03	0.59
26:1H:1532:C:H2'	26:1H:1533:C:O4'	2.03	0.59
19:AI:5:LEU:HB3	19:AI:10:PHE:HE2	1.68	0.59
1:1G:413:G:H2'	1:1G:428:G:N2	2.17	0.59
33:69:76:THR:HG23	33:69:77:LEU:H	1.67	0.59
3:22:40:ARG:HB3	3:22:55:VAL:HG11	1.84	0.59
26:14:1942:C:OP2	26:14:1943:U:O2'	2.13	0.59
4:3E:83:SER:HA	4:3E:89:THR:HG23	1.85	0.59
9:82:99:LEU:HB3	9:82:101:PHE:CE2	2.38	0.59
33:69:61:ARG:HA	33:69:64:GLU:HB2	1.85	0.59
38:55:20:LEU:HD21	38:55:40:LYS:HD3	1.84	0.59
1:13:1504:G:H3'	1:13:1504:G:P	2.43	0.59
26:14:573:G:O2'	26:14:574:C:H3'	2.03	0.59
1:13:1133:G:H2'	1:13:1134:G:C8	2.38	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:1H:1332:G:N2	26:1H:1610:A:C8	2.70	0.59
1:13:130:A:O2'	1:13:131:C:O5'	2.18	0.59
52:J5:3:LYS:O	52:J5:4:HIS:C	2.41	0.59
13:4A:15:VAL:HG12	13:4A:45:VAL:HG22	1.85	0.59
7:62:143:ARG:NH1	24:3L:41:C:O2'	2.35	0.59
17:8A:86:GLU:O	17:8A:90:ILE:HG12	2.03	0.59
2:12:70:PHE:HB2	2:12:92:TYR:HB2	1.84	0.59
1:13:598:U:H4'	8:7E:94:TYR:CG	2.38	0.59
26:14:1058:U:H2'	26:14:1059:G:H8	1.67	0.59
5:4E:33:VAL:HG11	5:4E:109:ILE:HA	1.84	0.59
43:E8:79:GLY:HA3	43:E8:100:THR:HG22	1.85	0.59
20:BI:26:ASN:O	20:BI:30:LYS:HB2	2.03	0.59
31:49:39:ILE:O	31:49:91:ARG:HG2	2.03	0.59
40:B8:80:SER:HB3	40:B8:83:ILE:HG13	1.84	0.59
1:13:1000:A:H2'	1:13:1001:G:C8	2.37	0.59
39:65:35:ILE:HB	39:65:97:ARG:NH2	2.18	0.59
47:I8:17:GLN:O	47:I8:19:LYS:HE3	2.02	0.59
43:E8:86:LEU:HD12	43:E8:87:PRO:HD2	1.85	0.59
26:14:2048:G:N7	58:14:3823:HOH:O	2.31	0.59
53:K5:27:LYS:O	53:K5:28:ARG:NH1	2.36	0.59
1:13:1218:C:H2'	1:13:1219:U:C6	2.38	0.59
37:45:134:ARG:O	37:45:136:ALA:N	2.35	0.59
26:14:1607:C:H4'	26:14:1608:A:O5'	2.03	0.59
2:12:7:VAL:HG22	2:12:8:LYS:H	1.68	0.58
26:1H:1774:C:OP1	58:1H:3745:HOH:O	2.16	0.58
16:7A:74:LEU:HD12	16:7A:79:VAL:HG21	1.85	0.58
26:1H:1021:A:H3'	26:1H:1022:G:H5''	1.84	0.58
51:I5:34:GLU:HG2	51:I5:35:VAL:H	1.68	0.58
17:8I:76:LEU:HD12	17:8I:77:VAL:H	1.68	0.58
35:25:10:VAL:HG21	35:25:16:ALA:O	2.03	0.58
15:6I:42:HIS:CD2	15:6I:43:LEU:HD23	2.38	0.58
1:13:1356:G:H2'	1:13:1357:A:C8	2.38	0.58
26:1H:2679:A:H4'	29:21:165:VAL:HG11	1.85	0.58
10:1A:99:LYS:HD3	10:1A:100:THR:H	1.67	0.58
28:11:201:HIS:O	28:11:204:ILE:HG12	2.03	0.58
26:1H:2734:A:H5'	26:1H:2735:G:OP2	2.03	0.58
26:14:1327:C:O3'	38:55:105:ARG:NH2	2.36	0.58
44:F8:3:THR:HA	44:F8:6:ASP:OD2	2.03	0.58
27:1J:104:A:H2'	27:1J:105:G:O4'	2.03	0.58
26:1H:1858:G:H2'	26:1H:1883:G:N2	2.17	0.58
26:14:289:A:H3'	26:14:290:G:H8	1.68	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
46:H8:99:TYR:HD1	46:H8:123:ASP:HB3	1.69	0.58
4:3E:170:VAL:HG22	4:3E:171:GLY:H	1.67	0.58
1:1G:823:G:H21	8:72:1:MET:HE2	1.68	0.58
52:J5:36:CYS:SG	52:J5:48:GLU:HB2	2.43	0.58
47:I8:24:LYS:O	47:I8:25:ARG:NH1	2.36	0.58
5:42:9:LYS:HG2	5:42:108:ALA:HB1	1.85	0.58
3:22:139:GLN:NE2	3:22:142:MET:SD	2.73	0.58
1:1G:804:U:H5'	1:1G:805:C:OP2	2.02	0.58
26:1H:1509:C:N3	26:1H:1511:A:N6	2.51	0.58
20:BI:30:LYS:NZ	20:BI:80:ARG:HH12	2.01	0.58
1:1G:1370:G:C8	9:82:109:VAL:HG11	2.38	0.58
26:1H:1537:C:H2'	26:1H:1538:G:C8	2.38	0.58
1:13:661:G:H1	1:13:744:C:H42	1.51	0.58
1:1G:501:C:H2'	1:1G:502:G:C8	2.36	0.58
26:14:646:A:H2'	26:14:647:G:O4'	2.03	0.58
9:8E:25:LYS:N	9:8E:60:ASP:OD1	2.35	0.58
26:14:579:G:H2'	26:14:580:C:C6	2.38	0.58
5:42:16:THR:OG1	5:42:16:THR:O	2.22	0.58
33:61:68:LEU:HA	33:61:71:ILE:HG22	1.85	0.58
49:K8:4:SER:HA	49:K8:6:VAL:HG22	1.85	0.58
26:1H:2179:C:H2'	26:1H:2180:U:C6	2.38	0.58
23:2K:62:C:H2'	23:2K:63:C:H6	1.68	0.58
4:32:55:ALA:O	4:32:59:ARG:HG2	2.03	0.58
26:14:860:U:O2'	26:14:861:A:H5'	2.02	0.58
27:16:15:A:H3'	27:16:16:G:H5'	1.84	0.58
40:B8:29:ARG:HB2	40:B8:46:GLU:HG3	1.86	0.58
9:8E:43:ALA:HA	9:8E:74:ILE:HD13	1.84	0.58
26:1H:1266:G:O4'	43:E8:15:ARG:NH2	2.35	0.58
1:13:1118:C:H1'	1:13:1179:A:C4	2.38	0.58
16:7I:39:TYR:OH	16:7I:41:PRO:HB3	2.03	0.58
26:1H:1432:C:H2'	26:1H:1433:U:O4'	2.03	0.58
28:19:37:LEU:HB2	28:19:38:LYS:HG2	1.84	0.58
26:1H:586:A:P	58:1H:3827:HOH:O	2.61	0.58
1:13:1346:A:H5''	9:8E:120:ARG:HH12	1.69	0.58
26:14:70:G:H21	26:14:71:A:H62	1.49	0.58
26:14:780:G:OP1	28:19:218:ARG:NH2	2.37	0.58
26:1H:598:G:C5'	36:78:11:GLY:HA3	2.32	0.58
31:49:122:PRO:O	31:49:125:PHE:HD2	1.87	0.58
26:1H:2315:G:OP1	31:41:36:LYS:NZ	2.33	0.58
26:14:527:C:H5	58:14:3922:HOH:O	1.85	0.58
30:31:114:VAL:HG21	30:31:202:PHE:CE2	2.39	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:25:88:ASN:HB3	35:25:94:ARG:HD3	1.86	0.58
1:1G:1157:A:H8	1:1G:1158:C:C5	2.22	0.58
1:13:1118:C:H1'	1:13:1179:A:C5	2.38	0.58
29:21:104:VAL:HG22	29:21:198:VAL:HG22	1.86	0.58
1:13:613:C:H42	1:13:627:G:H1	1.51	0.58
26:14:5:A:H2'	26:14:6:A:C8	2.39	0.58
43:E8:97:LYS:HE2	43:E8:99:ARG:NH2	2.18	0.58
26:1H:2492:U:H2'	26:1H:2493:U:C6	2.38	0.58
20:BI:50:GLU:HB2	20:BI:100:ILE:HB	1.84	0.58
1:13:221:C:H2'	1:13:222:U:H6	1.67	0.58
1:1G:987:G:N2	1:1G:1218:C:N3	2.51	0.58
42:95:58:VAL:HB	42:95:98:GLU:HB2	1.84	0.58
55:Q8:34:TRP:C	55:Q8:34:TRP:CD1	2.77	0.58
36:35:59:LEU:O	36:35:59:LEU:HD22	2.03	0.58
26:1H:1794:U:H2'	26:1H:1795:C:C6	2.39	0.58
1:13:711:G:OP1	6:5E:54:LYS:NZ	2.26	0.58
26:1H:1980:G:H4'	58:1H:3657:HOH:O	2.02	0.58
26:14:2786:U:H4'	29:29:64:LYS:C	2.23	0.58
25:4L:15:A:O5'	25:4L:15:A:H8	1.87	0.58
1:1G:520:A:OP1	12:3A:52:LEU:HB2	2.04	0.58
1:1G:1154:G:H2'	1:1G:1155:G:C8	2.32	0.58
40:B8:26:ASP:HB2	40:B8:91:ARG:HA	1.84	0.58
27:1J:40:U:H3	27:1J:43:C:H5''	1.69	0.58
1:13:872:A:C5	1:13:874:G:C8	2.91	0.58
29:29:33:VAL:HG12	29:29:89:ASP:HB3	1.85	0.58
1:1G:546:G:OP2	4:32:72:GLU:N	2.36	0.58
36:35:62:LEU:HD23	55:M5:27:THR:HG22	1.85	0.58
38:98:24:GLN:OE1	38:98:36:THR:HG21	2.03	0.58
26:14:2528:U:O3'	26:14:2529:G:N2	2.26	0.58
20:BI:53:LEU:HD12	20:BI:56:MET:HE2	1.85	0.58
1:1G:877:C:H5''	8:72:88:LYS:HD3	1.85	0.58
26:14:2062:A:O2'	26:14:2063:C:OP1	2.18	0.58
6:5E:76:ALA:O	6:5E:80:ARG:HG3	2.04	0.58
26:1H:1997:G:H5''	58:1H:3916:HOH:O	2.03	0.58
27:1J:46:A:H2'	27:1J:47:C:H6	1.69	0.58
26:14:654(D):G:H22	26:14:654(Q):C:N4	2.00	0.58
55:M5:48:PHE:HE2	55:M5:50:LEU:HD13	1.67	0.58
4:3E:36:ARG:HB3	4:3E:38:TYR:CZ	2.39	0.58
31:41:26:GLN:NE2	31:41:27:ASN:HB2	2.19	0.58
39:65:3:ARG:HE	39:65:4:LEU:N	2.01	0.58
26:14:2557:G:H2'	26:14:2558:C:C6	2.37	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:13:793:U:H5'	1:13:794:A:H5''	1.86	0.58
29:29:12:THR:HG22	40:75:58:ASN:HD21	1.69	0.58
33:69:101:LEU:H	33:69:101:LEU:HD23	1.67	0.58
44:B5:18:TYR:O	44:B5:20:GLY:N	2.36	0.58
48:J8:7:ILE:HG12	48:J8:62:VAL:HG11	1.86	0.58
7:6E:155:ARG:O	7:6E:155:ARG:NH2	2.35	0.58
26:1H:528:A:C2	26:1H:2043:C:H4'	2.38	0.58
1:1G:975:A:H4'	1:1G:976:G:H5''	1.85	0.58
46:D5:60:GLU:HA	46:D5:66:SER:HA	1.86	0.58
26:1H:2680:C:H5'	29:21:189:PRO:HA	1.86	0.58
1:1G:1302:U:OP1	13:4A:13:LYS:NZ	2.36	0.58
26:1H:1799:G:H5'	26:1H:1819:A:H61	1.68	0.58
27:1J:7:G:H1	27:1J:113:C:H42	1.52	0.58
28:11:17:THR:CG2	28:11:204:ILE:HA	2.34	0.58
26:14:2125:G:N2	26:14:2172:U:OP1	2.37	0.58
29:29:25:VAL:O	29:29:26:ILE:HG12	2.03	0.58
26:1H:2224:G:H4'	26:1H:2226:C:C2	2.39	0.58
1:13:247:G:OP2	17:8I:100:LYS:HB2	2.04	0.58
26:14:617:G:OP1	30:39:40:GLN:HG3	2.03	0.58
26:1H:1771:C:H1'	26:1H:1786:A:C8	2.39	0.58
26:1H:607:U:N3	26:1H:621:A:H2	1.94	0.58
54:L5:19:ARG:HG2	54:L5:19:ARG:HH11	1.69	0.58
1:1G:363:A:OP1	12:3A:33:ARG:HG3	2.03	0.58
37:45:117:ALA:HA	37:45:120:ILE:HB	1.85	0.58
1:13:343:U:O2'	1:13:346:G:O6	2.18	0.58
1:13:1240:U:C5	7:6E:32:ARG:HD2	2.37	0.58
1:13:1226:C:H4'	19:AI:80:TYR:OH	2.04	0.58
48:F5:85:LEU:HA	48:F5:87:PRO:HG2	1.85	0.58
44:F8:57:LEU:HD11	44:F8:78:LYS:HD2	1.86	0.58
26:1H:816:C:OP2	58:1H:3823:HOH:O	2.17	0.58
26:1H:729:G:O5'	28:11:208:LYS:NZ	2.37	0.58
26:1H:270(J):G:H2'	26:1H:270(K):C:O4'	2.04	0.58
5:42:140:ARG:O	5:42:143:ARG:NH2	2.37	0.58
1:13:116:A:H61	1:13:313:A:H1'	1.69	0.58
30:31:139:PHE:HB2	30:31:166:ALA:HB1	1.86	0.58
26:14:1161:C:H2'	26:14:1162:G:H8	1.69	0.58
1:13:195:A:H4'	20:BI:68:LYS:HE2	1.83	0.58
26:1H:1209:G:O6	58:1H:4267:HOH:O	2.16	0.58
37:45:66:ILE:HG13	37:45:67:ARG:H	1.69	0.58
32:51:97:ARG:NH2	32:51:104:GLU:OE2	2.33	0.58
26:1H:604:G:OP2	36:78:90:ARG:NH1	2.37	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:13:375:U:OP1	16:7I:69:THR:HG21	2.03	0.58
1:1G:537:G:H5''	12:3A:113:ARG:NH1	2.19	0.58
26:1H:654(D):G:N2	26:1H:654(R):C:N3	2.52	0.58
19:AI:39:THR:HG22	19:AI:40:ILE:H	1.68	0.58
26:14:2873:A:C8	38:55:5:LYS:HA	2.39	0.58
38:55:33:ARG:NH2	38:55:115:GLU:OE2	2.37	0.58
8:7E:87:SER:HB2	8:7E:93:VAL:CB	2.33	0.58
40:B8:108:ARG:HA	40:B8:111:ARG:NE	2.18	0.58
58:14:3700:HOH:O	30:39:55:GLY:HA2	2.04	0.58
4:3E:102:ASP:HB3	4:3E:136:PRO:HB2	1.85	0.58
26:1H:581:C:H2'	26:1H:582:G:H8	1.68	0.58
46:H8:7:ALA:HB2	46:H8:59:LEU:HD22	1.85	0.58
29:29:60:ASN:C	29:29:62:PRO:HD3	2.25	0.58
8:7E:8:ASP:O	8:7E:12:ARG:HG3	2.04	0.58
8:72:99:GLU:OE2	8:72:100:ILE:N	2.25	0.58
26:14:674:G:O2'	30:39:74:ARG:HG3	2.04	0.58
49:K8:35:LEU:HD12	49:K8:53:LEU:HD12	1.86	0.58
3:2E:95:THR:HB	3:2E:97:LYS:H	1.67	0.58
26:1H:1590:U:H2'	26:1H:1591:G:C8	2.38	0.58
26:14:605:C:O2	26:14:657:U:O2'	2.20	0.58
27:1J:88:C:H3'	27:1J:89:G:C8	2.39	0.58
26:1H:2311:A:C8	31:41:88:ILE:HD12	2.39	0.57
32:59:163:TYR:CE1	32:59:169:VAL:HG21	2.39	0.57
25:4L:12:A:H4'	25:4L:13:A:OP2	2.04	0.57
12:3A:62:SER:HB2	12:3A:64:TYR:CD1	2.38	0.57
26:1H:2261:C:H1'	26:1H:2388:A:N3	2.19	0.57
3:2E:7:PRO:O	3:2E:11:ARG:HG2	2.04	0.57
22:1K:52:G:H5'	37:88:56:ARG:HH21	1.68	0.57
39:65:34:HIS:CD2	39:65:54:LEU:HD22	2.39	0.57
10:1I:65:LEU:HD13	14:5I:56:VAL:HG22	1.86	0.57
48:J8:18:ILE:HG12	48:J8:37:ILE:HG12	1.85	0.57
10:1A:24:VAL:HG13	10:1A:34:VAL:HG11	1.86	0.57
41:C8:32:PHE:HZ	41:C8:36:ARG:HH21	1.52	0.57
34:58:15:LEU:HB2	34:58:134:ARG:HB3	1.85	0.57
26:1H:825:C:H5''	58:1H:3765:HOH:O	2.02	0.57
26:14:2656:U:H3	26:14:2665:A:H2	1.52	0.57
31:41:20:ILE:HG23	31:41:25:TYR:HB2	1.86	0.57
26:1H:1210:A:OP1	26:1H:1211:U:O2'	2.21	0.57
26:14:72:U:OP2	49:G5:29:LYS:NZ	2.30	0.57
1:13:1160:G:H1	1:13:1177:G:H1	1.52	0.57
27:1J:46:A:H2'	27:1J:47:C:C6	2.39	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:13:686:U:O4	1:13:703:G:H1'	2.04	0.57
26:14:2015:A:H1'	52:J5:2:ALA:HA	1.86	0.57
26:1H:2400:G:H2'	26:1H:2401:U:H6	1.69	0.57
46:D5:128:VAL:HG22	46:D5:129:SER:H	1.69	0.57
46:H8:77:ASP:OD2	46:H8:80:ARG:NH1	2.38	0.57
32:51:126:PRO:HG2	32:51:130:ARG:HH12	1.68	0.57
2:1E:124:SER:HB2	2:1E:125:PRO:HD2	1.86	0.57
4:32:24:GLU:HG2	4:32:25:ARG:H	1.70	0.57
1:13:1178:G:N7	9:8E:97:LYS:NZ	2.52	0.57
1:13:1004:A:H1'	1:13:1036:G:H22	1.68	0.57
32:51:4:ILE:HG21	32:51:6:ARG:NH1	2.18	0.57
29:21:111:ARG:HD3	29:21:160:TYR:CE2	2.39	0.57
26:1H:1331:A:O2'	26:1H:1332:G:H8	1.86	0.57
26:1H:2345:G:H4'	26:1H:2346:A:O5'	2.03	0.57
1:1G:1179:A:H4'	9:82:103:THR:HA	1.86	0.57
1:1G:1141:C:H2'	1:1G:1142:G:C8	2.39	0.57
26:14:2064:C:H2'	26:14:2065:C:C6	2.40	0.57
1:13:455:C:H42	1:13:477:G:H1	1.51	0.57
29:21:29:GLY:H	29:21:51:PHE:HE1	1.49	0.57
26:14:2232:U:P	48:F5:40:ARG:HH12	2.27	0.57
1:13:1051:C:H2'	1:13:1052:U:H6	1.69	0.57
4:3E:94:LEU:HA	4:3E:97:LEU:HD12	1.85	0.57
37:45:26:TYR:O	37:45:28:ALA:N	2.36	0.57
34:58:73:THR:HB	34:58:82:LEU:HD11	1.86	0.57
54:P8:5:TRP:NE1	54:P8:7:PRO:HG3	2.19	0.57
1:1G:1360:A:OP1	1:1G:1360:A:H8	1.88	0.57
31:41:5:VAL:H	51:M8:25:TYR:HE2	1.52	0.57
26:1H:1702:G:O6	58:1H:4087:HOH:O	2.14	0.57
26:14:2392:A:H2	26:14:2424:C:N4	2.00	0.57
41:C8:92:ARG:HD2	42:D8:11:GLN:HB2	1.86	0.57
37:45:81:VAL:O	37:45:82:ARG:NH1	2.37	0.57
26:1H:2298:A:H2'	26:1H:2299:G:O4'	2.04	0.57
26:1H:2684:U:C4	26:1H:2685:G:N7	2.72	0.57
26:1H:1602:U:O4	58:1H:4224:HOH:O	2.18	0.57
29:21:64:LYS:O	29:21:70:ALA:HB2	2.05	0.57
7:6E:73:MET:HG2	7:6E:90:GLU:HA	1.85	0.57
2:12:19:HIS:CD2	2:12:20:GLU:HG2	2.39	0.57
24:3K:10:G:H1	24:3K:25:C:H42	1.52	0.57
26:14:1292:U:H2'	26:14:1293:C:C6	2.39	0.57
24:3L:18:G:H2'	24:3L:57:G:H22	1.69	0.57
26:14:945:A:P	58:14:3658:HOH:O	2.61	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
27:16:94:C:H2'	27:16:95:U:C6	2.40	0.57
40:75:45:PHE:CE2	40:75:74:ARG:HD3	2.39	0.57
2:1E:97:TRP:CZ3	2:1E:172:ILE:HB	2.40	0.57
27:1J:75:G:H21	46:D5:85:HIS:CE1	2.22	0.57
30:31:164:ARG:HH11	30:31:164:ARG:HG2	1.68	0.57
14:5A:12:ARG:HD3	14:5A:13:THR:H	1.69	0.57
8:72:27:PRO:O	8:72:32:LYS:HD3	2.04	0.57
26:14:363(E):U:H5'	26:14:363(F):A:OP2	2.04	0.57
55:Q8:49:VAL:HG13	55:Q8:50:LEU:N	2.18	0.57
35:68:68:GLU:OE2	35:68:78:ARG:NH1	2.37	0.57
1:1G:1131:G:H2'	1:1G:1132:C:H6	1.68	0.57
18:9A:22:VAL:HG22	18:9A:23:LYS:H	1.68	0.57
1:1G:1226:C:H3'	13:4A:96:LEU:HD21	1.86	0.57
33:69:101:LEU:O	33:69:106:GLY:N	2.38	0.57
26:14:855:G:H5''	26:14:856:C:OP2	2.05	0.57
1:1G:668:G:H4'	15:6A:48:LYS:HB3	1.85	0.57
26:1H:2689:U:H5''	26:1H:2713:A:C2	2.40	0.57
4:3E:175:SER:HB3	4:3E:186:LEU:HD21	1.86	0.57
26:1H:2111:C:H2'	26:1H:2118:U:H4'	1.86	0.57
26:1H:2147:G:H2'	26:1H:2148:G:H4'	1.85	0.57
1:1G:330:C:O2	58:1G:1713:HOH:O	2.17	0.57
33:61:56:LYS:O	33:61:60:GLU:HB3	2.04	0.57
36:35:78:PRO:HB3	36:35:111:ARG:NH2	2.19	0.57
1:13:280:C:H3'	1:13:281:G:H5'	1.85	0.57
1:13:664:G:N2	1:13:741:G:H1	1.97	0.57
41:85:92:ARG:NH1	42:95:11:GLN:H	2.01	0.57
26:1H:654(D):G:H1	26:1H:654(Q):C:H42	1.50	0.57
1:1G:957:U:H1'	1:1G:960:U:H5	1.70	0.57
1:13:973:G:H3'	1:13:974:A:H5''	1.87	0.57
24:3K:7:A:O2'	24:3K:49:C:H5'	2.05	0.57
41:C8:92:ARG:CZ	42:D8:11:GLN:H	2.18	0.57
26:14:619:G:H5''	26:14:620:G:N2	2.20	0.57
26:1H:2287:A:N6	26:1H:2344:U:H3	2.03	0.57
26:1H:1278:A:H4'	38:98:34:ILE:HD11	1.87	0.57
7:62:113:GLU:O	7:62:119:ARG:HD3	2.05	0.57
1:13:1286:A:H5''	21:1F:26:LYS:HG2	1.86	0.57
26:1H:2469:A:H61	26:1H:2481:G:H1'	1.70	0.57
6:52:61:LEU:HD23	6:52:63:TYR:OH	2.04	0.57
3:2E:128:PHE:CZ	3:2E:132:ARG:HD2	2.40	0.57
26:14:1288:U:C2	26:14:1327:C:O2	2.58	0.57
53:O8:41:PRO:HB2	53:O8:43:CYS:H	1.69	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:2688:U:H5	26:14:2720:U:OP2	1.88	0.57
10:1A:51:ARG:HE	10:1A:61:GLU:HB2	1.69	0.57
27:16:80:U:H2'	27:16:81:G:H21	1.69	0.57
26:1H:1374:G:H2'	26:1H:1375:C:C6	2.39	0.57
29:29:12:THR:HG22	40:75:58:ASN:ND2	2.19	0.57
1:13:673:G:H2'	1:13:674:G:C8	2.39	0.57
26:14:2111:C:N4	26:14:2147:G:H21	2.03	0.57
35:25:43:VAL:HG23	35:25:56:ASP:O	2.04	0.57
46:H8:62:PRO:C	46:H8:64:GLY:HA2	2.24	0.57
1:13:1342:C:O2'	9:8E:124:GLN:HG2	2.04	0.57
26:1H:996:A:H4'	41:C8:92:ARG:HE	1.68	0.57
1:13:1003:G:N2	1:13:1004:A:O2'	2.37	0.57
1:1G:192:U:H2'	1:1G:193:C:C6	2.35	0.57
26:14:2438:U:O3'	26:14:2439:A:H3'	2.05	0.57
1:13:411:A:C4	1:13:413:G:H1'	2.39	0.57
34:15:56:ASN:H	34:15:125:GLY:HA3	1.68	0.57
41:C8:17:ILE:HD12	41:C8:32:PHE:CE1	2.40	0.57
41:C8:50:ARG:HH12	42:D8:72:VAL:HG12	1.69	0.57
28:11:46:GLN:HB2	28:11:48:ARG:HG2	1.87	0.57
30:39:110:LEU:HD13	30:39:205:ARG:HG2	1.86	0.57
5:4E:76:ILE:HG13	5:4E:93:PRO:HB3	1.87	0.57
1:13:21:G:OP1	58:13:1814:HOH:O	2.17	0.57
26:14:300:A:N6	58:14:3887:HOH:O	2.38	0.57
39:65:87:PHE:CZ	39:65:102:ALA:HB2	2.39	0.57
8:72:18:ARG:NH2	8:72:81:HIS:O	2.38	0.57
26:1H:99:U:O4	45:G8:8:LYS:NZ	2.31	0.57
40:75:92:GLY:HA2	40:75:116:ALA:HA	1.86	0.57
38:98:42:LYS:HA	38:98:45:ARG:HD2	1.87	0.57
6:52:81:ILE:HD11	28:19:125:ILE:HG12	1.85	0.57
11:2I:19:ALA:O	11:2I:82:VAL:HA	2.04	0.57
1:13:31:G:O2'	1:13:48:C:N4	2.37	0.57
2:12:8:LYS:HE2	2:12:213:LEU:HD21	1.87	0.57
26:14:531:C:OP1	26:14:561:G:C2	2.57	0.57
41:85:92:ARG:NH2	42:95:10:LYS:HA	2.19	0.57
24:3K:66:U:H2'	24:3K:67:C:C6	2.40	0.57
26:14:2420:C:H41	55:M5:31:HIS:HB3	1.70	0.57
26:1H:1550:C:H2'	26:1H:1551:C:H6	1.70	0.57
1:1G:1372:U:OP2	9:82:11:LYS:NZ	2.24	0.57
26:14:2207:C:H42	26:14:2217:G:H1	1.52	0.57
31:41:35:GLU:OE1	31:41:36:LYS:N	2.37	0.57
1:13:1497:G:C2'	1:13:1498:U:H5'	2.35	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:981:A:N1	26:14:2027:G:O2'	2.27	0.57
26:14:706:A:H2'	26:14:707:G:O4'	2.04	0.57
47:E5:23:VAL:HG13	47:E5:38:VAL:HG22	1.87	0.57
26:1H:299:A:H5'	26:1H:300:A:OP2	2.04	0.57
26:14:698:C:O2'	26:14:734:A:N6	2.35	0.57
35:25:68:GLU:HB3	35:25:78:ARG:HB2	1.85	0.57
23:2K:22:A:C8	23:2K:22:A:H5''	2.40	0.57
33:69:63:ALA:HA	33:69:66:GLU:HG2	1.86	0.57
45:C5:29:GLU:N	45:C5:29:GLU:OE1	2.38	0.57
39:65:25:ARG:HD3	39:65:40:ILE:HD12	1.87	0.57
29:29:119:ARG:HG3	29:29:160:TYR:HB2	1.86	0.57
9:82:102:LEU:O	9:82:103:THR:OG1	2.20	0.57
1:1G:1326:C:H2'	1:1G:1327:C:C6	2.39	0.57
12:3I:47:LYS:HA	12:3I:49:ASN:H	1.70	0.57
22:1K:15:G:N2	22:1K:59:U:O2	2.36	0.57
2:1E:141:GLU:O	2:1E:145:LEU:HB2	2.05	0.57
1:13:244:U:H4'	1:13:245:C:O5'	2.05	0.57
1:13:56:U:H4'	33:69:82:ARG:HH21	1.69	0.57
35:68:86:ILE:HG22	35:68:94:ARG:HG3	1.86	0.57
1:13:1289:A:H3'	1:13:1290:G:H8	1.69	0.57
20:BA:11:SER:HA	20:BA:13:LEU:H	1.68	0.57
27:1J:15:A:H3'	27:1J:16:G:H5'	1.86	0.56
26:1H:511:U:C5	26:1H:512:G:C5	2.93	0.56
29:21:146:THR:HA	29:21:147:PRO:C	2.25	0.56
1:13:1157:A:N6	1:13:1178:G:N3	2.53	0.56
26:14:2210:G:O5'	26:14:2211:G:N2	2.38	0.56
1:1G:983:A:H2	1:1G:984:C:C6	2.23	0.56
30:31:65:TRP:CZ3	30:31:72:ARG:HB3	2.40	0.56
26:1H:960:A:H61	37:88:82:ARG:NH2	2.04	0.56
2:1E:87:ARG:NH1	2:1E:233:SER:HB2	2.19	0.56
23:2L:9:G:O2'	23:2L:10:G:N7	2.24	0.56
4:32:173:TRP:CD1	4:32:174:LEU:HG	2.39	0.56
50:L8:10:LYS:HD3	50:L8:53:LEU:HD23	1.87	0.56
1:13:624:C:H2'	1:13:625:G:C8	2.39	0.56
3:22:11:ARG:HB2	3:22:11:ARG:HH11	1.69	0.56
46:H8:63:ASP:N	46:H8:64:GLY:HA2	2.20	0.56
12:3I:82:VAL:HG12	12:3I:106:ASP:OD2	2.05	0.56
26:14:1335:U:OP1	44:B5:65:ARG:NH1	2.38	0.56
2:12:77:ALA:HB2	2:12:211:ILE:HD13	1.86	0.56
26:14:2464:C:H2'	26:14:2465:C:O4'	2.04	0.56
4:3E:106:TYR:HE2	4:3E:107:ARG:HH11	1.53	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:13:179:A:H2'	1:13:180:U:H6	1.69	0.56
1:1G:448:A:H2'	1:1G:449:C:O2	2.05	0.56
26:14:863:A:H2	26:14:914:C:H41	1.51	0.56
24:1L:10:G:H2'	24:1L:11:C:C6	2.40	0.56
1:1G:562:C:H1'	12:3A:15:ARG:HD2	1.87	0.56
47:E5:64:ASP:OD1	47:E5:64:ASP:N	2.38	0.56
20:BI:14:LYS:HG3	20:BI:17:ARG:HE	1.70	0.56
51:M8:41:PRO:HA	51:M8:47:GLN:HG3	1.87	0.56
5:4E:126:ARG:NH1	5:4E:126:ARG:HG3	2.17	0.56
26:14:2312:U:OP1	31:49:74:LYS:HD2	2.06	0.56
47:E5:68:GLU:OE2	47:E5:82:ARG:NH2	2.38	0.56
19:AI:9:VAL:HG21	51:M8:63:TYR:O	2.04	0.56
32:59:6:ARG:H	32:59:6:ARG:HD2	1.70	0.56
32:51:83:TYR:HB3	32:51:134:SER:HA	1.87	0.56
1:1G:1320:C:N3	19:AA:36:ARG:NH2	2.54	0.56
26:14:1328:G:H2'	26:14:1330:C:C5	2.40	0.56
13:4I:11:ARG:HG3	13:4I:12:ASN:N	2.20	0.56
26:14:1537:C:H2'	26:14:1538:G:C8	2.39	0.56
31:41:102:PHE:HA	31:41:105:LYS:HE2	1.87	0.56
40:75:45:PHE:CE2	40:75:74:ARG:HB2	2.40	0.56
11:2A:21:ILE:HB	11:2A:84:VAL:HG12	1.86	0.56
26:1H:1381:G:N7	58:1H:3924:HOH:O	2.33	0.56
9:8E:125:TYR:HD1	9:8E:126:SER:H	1.53	0.56
31:49:97:ASP:HA	31:49:100:TRP:HD1	1.70	0.56
36:35:52:GLU:O	36:35:54:GLY:N	2.38	0.56
26:1H:2376:A:H2	39:A8:112:PHE:HB3	1.70	0.56
26:1H:2789:C:H1'	26:1H:2892:A:H2	1.69	0.56
30:39:101:LEU:O	30:39:106:ARG:NH1	2.38	0.56
1:13:277:C:H2'	1:13:278:G:H8	1.71	0.56
3:2E:15:THR:HG21	3:2E:181:ASN:HA	1.87	0.56
26:14:2404:C:O3'	36:35:77:ARG:NH2	2.38	0.56
26:14:1839:G:C8	26:14:1927:A:H1'	2.41	0.56
29:29:55:ASN:O	29:29:57:LYS:N	2.33	0.56
2:1E:5:ILE:HG13	2:1E:6:THR:H	1.70	0.56
1:1G:540:G:H2'	1:1G:541:G:O4'	2.05	0.56
26:14:2318:G:H5'	26:14:2319:G:OP2	2.05	0.56
15:6I:6:GLU:H	15:6I:6:GLU:CD	2.09	0.56
1:1G:373:A:H2'	1:1G:374:A:H8	1.70	0.56
1:13:201:C:N4	1:13:209:U:H1'	2.21	0.56
27:16:15:A:H5'	27:16:16:G:C8	2.40	0.56
26:1H:833:U:O2	36:78:55:ARG:NH2	2.38	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
39:65:34:HIS:NE2	39:65:54:LEU:HD13	2.20	0.56
1:1G:560:U:O2'	1:1G:561:U:OP2	2.21	0.56
24:3L:52:G:N2	24:3L:63:G:N7	2.53	0.56
4:3E:65:ARG:NH1	4:3E:70:ILE:O	2.32	0.56
26:1H:2257:U:O2'	26:1H:2258:C:H5'	2.05	0.56
7:62:115:ARG:HB2	7:62:118:VAL:HG13	1.87	0.56
26:1H:363(B):G:H2'	26:1H:363(C):G:H8	1.70	0.56
36:35:11:GLY:C	36:35:13:ASN:H	2.06	0.56
1:1G:600:C:H2'	1:1G:601:C:C6	2.39	0.56
26:1H:1486:A:H2'	26:1H:1487:G:H8	1.71	0.56
26:14:1224:G:N2	26:14:1227:A:OP2	2.33	0.56
7:6E:11:GLN:OE1	7:6E:12:LEU:N	2.39	0.56
4:3E:141:ARG:HB2	4:3E:141:ARG:NH1	2.20	0.56
26:1H:236:C:H2'	26:1H:237:C:H6	1.70	0.56
39:A8:10:ARG:O	39:A8:14:VAL:HG13	2.05	0.56
28:11:6:PHE:HE2	28:11:18:VAL:HG23	1.71	0.56
26:1H:2098:U:H3	26:1H:2191:G:H1	1.54	0.56
1:13:1194:U:H2'	1:13:1195:C:C6	2.39	0.56
20:BI:69:GLY:O	20:BI:73:HIS:NE2	2.38	0.56
2:1E:189:ASP:HB3	2:1E:191:ASP:HB2	1.87	0.56
1:13:600:C:H2'	1:13:601:C:H6	1.71	0.56
55:Q8:35:GLN:C	55:Q8:37:SER:H	2.08	0.56
39:65:3:ARG:HE	39:65:4:LEU:H	1.53	0.56
46:D5:128:VAL:HG11	46:D5:133:ILE:HG23	1.88	0.56
45:C5:52:SER:HA	45:C5:55:TYR:O	2.05	0.56
17:8I:13:ASP:HA	17:8I:19:VAL:HG12	1.87	0.56
26:1H:107:C:H2'	26:1H:108:U:H6	1.70	0.56
33:69:38:LEU:HD12	33:69:38:LEU:H	1.71	0.56
1:13:814:A:N7	1:13:816:A:C4	2.73	0.56
26:1H:1336:A:H2'	26:1H:1337:G:C8	2.40	0.56
46:D5:124:ILE:HD11	46:D5:165:VAL:HG11	1.87	0.56
45:G8:87:LYS:HD2	45:G8:96:ILE:HD11	1.87	0.56
1:1G:519:C:H2'	1:1G:520:A:O4'	2.04	0.56
55:Q8:53:PRO:HA	55:Q8:54:GLU:C	2.26	0.56
26:14:328:U:H4'	45:C5:68:HIS:CD2	2.41	0.56
43:A5:11:ARG:HD2	43:A5:82:LEU:HD12	1.86	0.56
1:1G:1286:A:H3'	1:1G:1286:A:C8	2.40	0.56
26:1H:2781:A:C5'	26:1H:2782:G:H5'	2.36	0.56
26:1H:1103:A:H3'	26:1H:1104:C:H6	1.70	0.56
1:13:1247:U:H3	1:13:1290:G:H1	1.54	0.56
8:7E:127:LEU:HB2	8:7E:129:VAL:HG22	1.86	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:1217:C:OP1	41:85:15:LYS:NZ	2.29	0.56
26:1H:2233:U:H2'	26:1H:2234:G:C8	2.40	0.56
29:29:182:LEU:O	29:29:183:LEU:HD12	2.06	0.56
32:51:86:GLU:H	32:51:86:GLU:CD	2.07	0.56
26:14:800:A:OP1	58:14:3856:HOH:O	2.17	0.56
2:12:73:THR:HG21	2:12:97:TRP:N	2.19	0.56
26:14:2377:A:H4'	39:65:111:GLU:HG2	1.86	0.56
42:95:48:GLY:HA3	42:95:51:VAL:C	2.26	0.56
26:14:639:U:H2'	26:14:640:C:C6	2.40	0.56
26:14:1784:A:H4'	26:14:1785:A:O5'	2.06	0.56
37:45:25:ASP:OD1	37:45:25:ASP:N	2.37	0.56
26:1H:582:G:H1	26:1H:1258:C:H42	1.54	0.56
3:22:21:ARG:NH1	3:22:21:ARG:HB3	2.21	0.56
34:58:13:TRP:O	34:58:134:ARG:HA	2.06	0.56
1:1G:1080:A:H5'	5:42:14:ARG:NH2	2.20	0.56
28:19:228:PRO:HD3	28:19:234:GLY:O	2.06	0.56
26:14:1386:C:H2'	26:14:1387:C:C6	2.40	0.56
35:25:76:ALA:HB3	40:75:75:ILE:HB	1.88	0.56
26:14:2432:A:C8	48:F5:33:LYS:HD2	2.40	0.56
11:2I:41:THR:HG22	11:2I:42:TRP:H	1.70	0.56
7:62:15:ASP:OD1	7:62:44:TYR:OH	2.22	0.56
26:1H:1827:C:C2'	26:1H:1828:G:H5'	2.35	0.56
2:12:91:PRO:HG2	2:12:155:LEU:HB2	1.88	0.56
1:1G:957:U:H1'	1:1G:960:U:C5	2.40	0.56
1:1G:376:G:O3'	16:7A:5:ARG:NH1	2.35	0.56
26:14:67:U:H2'	26:14:68:G:H8	1.71	0.56
31:41:67:LYS:HE2	31:41:67:LYS:H	1.71	0.56
26:1H:1021:A:H8	26:1H:1021:A:H3'	1.70	0.56
1:1G:1028(A):C:H42	1:1G:1032(B):G:H22	1.52	0.56
32:51:157:TYR:CE1	32:51:172:LYS:HB2	2.40	0.56
26:1H:1709:U:H1'	26:1H:2860:A:N3	2.20	0.56
13:4I:27:LYS:HD3	13:4I:31:LYS:HZ3	1.70	0.56
26:1H:1433:U:O2	26:1H:1561:G:C2	2.58	0.56
36:35:78:PRO:HA	36:35:110:TYR:CE2	2.41	0.56
26:1H:130:C:O3'	26:1H:1349:A:H1'	2.05	0.56
31:49:37:VAL:HG22	31:49:159:VAL:HG12	1.87	0.56
1:1G:297:G:N2	1:1G:300:A:OP2	2.38	0.56
29:29:147:PRO:HB2	29:29:149:ARG:HG2	1.88	0.56
26:14:185:U:H4'	26:14:218:A:H4'	1.88	0.56
1:1G:313:A:H2'	1:1G:314:C:C6	2.41	0.56
41:C8:68:ALA:O	41:C8:71:GLN:HB2	2.06	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:2880:C:H1'	38:55:92:GLY:HA3	1.86	0.56
1:13:426:G:OP1	4:3E:38:TYR:OH	2.22	0.56
36:35:63:PRO:HB3	55:M5:13:ARG:HG3	1.87	0.56
23:2L:41:C:H2'	23:2L:42:C:C6	2.39	0.56
26:14:2529:G:P	26:14:2529:G:H21	2.28	0.56
26:1H:1266:G:O2'	26:1H:2012:G:O6	2.18	0.56
22:1K:51:U:H2'	22:1K:52:G:H8	1.71	0.56
31:41:17:PRO:HA	31:41:20:ILE:HG13	1.88	0.56
37:45:26:TYR:O	37:45:26:TYR:HD1	1.87	0.56
27:16:94:C:H2'	27:16:95:U:H6	1.70	0.56
40:75:45:PHE:CD2	40:75:74:ARG:HD3	2.40	0.56
26:14:1461:G:H2'	26:14:1462:C:C6	2.41	0.56
9:82:24:GLY:HA2	9:82:59:PHE:O	2.06	0.56
34:15:1:MET:O	34:15:2:LYS:HG2	2.06	0.56
42:D8:10:LYS:NZ	42:D8:23:GLU:OE1	2.38	0.56
26:14:2126:A:N1	26:14:2163:C:H1'	2.20	0.56
41:C8:112:ARG:CZ	42:D8:47:VAL:HG22	2.36	0.56
33:69:112:LYS:HA	33:69:114:LEU:H	1.70	0.56
26:1H:259:G:C2'	26:1H:621:A:HO2'	2.18	0.56
30:31:101:LEU:O	30:31:106:ARG:NH1	2.38	0.56
27:16:40:U:H1'	27:16:45:A:N6	2.21	0.56
26:1H:1021:A:C8	26:1H:1021:A:H3'	2.39	0.56
26:1H:1657:C:H2'	26:1H:1658:C:C6	2.40	0.56
45:G8:97:ARG:HB2	45:G8:102:CYS:HB2	1.88	0.56
26:1H:460:A:H5''	26:1H:461:C:OP2	2.06	0.56
26:14:2537:U:H2'	26:14:2538:C:H6	1.68	0.56
26:1H:1590:U:H2'	26:1H:1591:G:H8	1.69	0.56
8:72:17:THR:HB	8:72:18:ARG:HH11	1.70	0.56
28:11:111:LEU:HD23	28:11:127:VAL:HG12	1.88	0.56
26:1H:817:C:H4'	26:1H:932:G:C5	2.40	0.56
26:1H:2123:G:H2'	26:1H:2124:G:H8	1.71	0.56
30:39:135:LYS:HB3	30:39:138:GLU:HG3	1.88	0.56
26:1H:2331:G:O3'	47:I8:43:THR:HG22	2.05	0.56
10:1I:29:ARG:HH22	10:1I:84:GLN:HE22	1.52	0.56
1:1G:793:U:H5'	1:1G:794:A:O5'	2.06	0.56
23:2L:13:C:O2'	26:14:1924:C:H4'	2.05	0.56
26:14:531:C:H4'	26:14:532:A:H5''	1.87	0.56
26:1H:2032:G:H1'	29:21:145:LYS:HE3	1.87	0.56
26:1H:1556:C:H2'	26:1H:1557:C:H6	1.71	0.56
4:32:31:CYS:C	4:32:33:MET:H	2.06	0.56
31:41:40:ASN:HB2	31:41:91:ARG:HG2	1.88	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:95:G:H4'	49:G5:46:GLN:HB2	1.88	0.56
2:1E:70:PHE:HE2	2:1E:90:MET:HB3	1.71	0.56
26:14:1871:A:H2'	26:14:1872:A:C8	2.40	0.56
1:13:598:U:H4'	8:7E:94:TYR:CD2	2.41	0.56
39:65:62:LYS:O	39:65:66:ALA:N	2.39	0.56
34:58:57:ALA:C	34:58:59:LYS:H	2.08	0.56
13:4I:82:MET:HG2	13:4I:93:ARG:HG3	1.88	0.56
26:14:597:U:H2'	26:14:598:G:C8	2.41	0.56
26:14:606:U:H4'	26:14:658:C:H4'	1.88	0.56
26:1H:859:G:H5'	26:1H:2268:A:O2'	2.06	0.56
1:13:1096:C:H2'	1:13:1097:C:H6	1.70	0.56
1:1G:522:C:OP2	12:3A:69:TYR:OH	2.24	0.56
1:13:1378:C:O2	7:6E:76:ARG:NH1	2.39	0.56
26:1H:637:A:H2'	36:78:117:GLU:OE1	2.06	0.56
1:13:1130:A:O2'	9:8E:3:GLN:NE2	2.32	0.55
23:2L:8:4SU:C2	23:2L:14:A:H62	2.19	0.55
26:1H:1063:G:N2	26:1H:1076:C:O2'	2.39	0.55
26:14:890:A:H2'	26:14:892:G:H8	1.71	0.55
1:1G:1294:G:H2'	1:1G:1295:G:C8	2.42	0.55
13:4A:57:ARG:HH12	51:I5:17:GLY:HA3	1.70	0.55
26:1H:1711:C:H2'	26:1H:1712:C:C6	2.42	0.55
26:1H:1337:G:H2'	26:1H:1338:G:H8	1.70	0.55
26:1H:806:C:C2	26:1H:807:U:C5	2.94	0.55
26:1H:2292:C:OP1	39:A8:17:ARG:NH2	2.34	0.55
20:BI:35:THR:HA	20:BI:38:LYS:HD3	1.88	0.55
26:14:1053:C:H42	26:14:1106:G:H1	1.54	0.55
26:1H:1568:G:H5''	28:11:61:LEU:HD22	1.86	0.55
10:1A:33:GLN:HB3	10:1A:75:ILE:HG12	1.87	0.55
15:6A:39:LEU:HD12	15:6A:56:LEU:HD13	1.88	0.55
2:1E:236:TYR:HA	2:1E:239:VAL:HG21	1.87	0.55
26:14:2793:G:N2	26:14:2804:C:N3	2.54	0.55
39:A8:100:ALA:HA	39:A8:103:GLU:HG2	1.88	0.55
26:1H:1963:U:O5'	26:1H:1963:U:H6	1.88	0.55
26:1H:234:C:H2'	26:1H:235:U:H6	1.71	0.55
36:78:19:VAL:HB	36:78:20:GLY:HA2	1.88	0.55
26:1H:1899:G:N2	26:1H:1902:C:C5	2.74	0.55
26:14:1204:A:O2'	26:14:1205:U:OP2	2.23	0.55
36:35:39:LYS:CD	36:35:45:LEU:HD21	2.34	0.55
26:14:996:A:H1'	41:85:92:ARG:NH2	2.22	0.55
26:14:996:A:H2'	26:14:997:G:H8	1.71	0.55
42:95:71:LEU:CA	42:95:86:GLY:HA2	2.36	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
53:K5:25:LYS:HB2	55:M5:34:TRP:CZ3	2.41	0.55
26:14:1557:C:H5''	26:14:1558:A:OP2	2.07	0.55
1:1G:707:C:H2'	1:1G:708:C:C6	2.41	0.55
4:32:61:LYS:NZ	4:32:72:GLU:OE2	2.37	0.55
11:2A:43:SER:OG	11:2A:44:SER:N	2.38	0.55
1:1G:861:G:H2'	1:1G:862:C:C6	2.41	0.55
26:1H:1692:U:O2'	26:1H:1693:U:H2'	2.06	0.55
53:K5:28:ARG:HG3	53:K5:31:PRO:HD2	1.88	0.55
26:1H:1448:G:H1'	26:1H:1528:A:H62	1.72	0.55
29:29:94:GLU:HG3	29:29:177:PRO:HB3	1.88	0.55
29:21:46:ALA:HA	29:21:82:ARG:O	2.05	0.55
32:51:37:VAL:HG22	32:51:68:THR:HG23	1.88	0.55
41:85:52:ARG:HA	41:85:55:ARG:HD3	1.89	0.55
28:11:12:SER:O	28:11:16:MET:HB2	2.06	0.55
30:31:179:GLU:OE1	30:31:179:GLU:N	2.38	0.55
26:14:2562:U:O3'	35:25:25:LEU:HD21	2.06	0.55
12:3A:47:LYS:HG2	12:3A:48:PRO:HD2	1.88	0.55
1:1G:963:G:N2	10:1A:55:LYS:HE3	2.07	0.55
26:1H:1359:A:C2	26:1H:1372:U:O4	2.59	0.55
26:1H:2392:A:H2	26:1H:2424:C:N4	2.00	0.55
26:1H:1265:A:OP1	26:1H:1265:A:H8	1.89	0.55
26:1H:1517:G:H5''	26:1H:1518:C:OP2	2.05	0.55
1:13:963:G:N2	10:1I:55:LYS:HZ1	2.04	0.55
26:14:2392:A:H8	36:35:61:ARG:HD2	1.71	0.55
1:13:376:G:O3'	16:7I:5:ARG:NH1	2.26	0.55
1:13:1149:C:H2'	1:13:1150:U:C6	2.40	0.55
40:75:4:GLY:HA2	40:75:8:LYS:HZ2	1.71	0.55
1:1G:1171:G:H2'	1:1G:1172:C:C6	2.42	0.55
1:1G:1109:C:H2'	1:1G:1110:A:O4'	2.07	0.55
35:25:104:ARG:NH1	40:75:36:GLU:HB3	2.21	0.55
26:1H:67:U:H3	26:1H:74:A:H2	1.55	0.55
2:1E:115:LEU:HD13	2:1E:145:LEU:HB3	1.88	0.55
42:D8:47:VAL:HG23	42:D8:48:GLY:N	2.21	0.55
1:13:50:A:H1'	1:13:52:G:C8	2.41	0.55
26:1H:1168:G:C2	26:1H:1182:A:C2	2.94	0.55
1:13:517:G:N1	1:13:533:A:OP2	2.35	0.55
1:1G:1310:G:OP1	13:4A:77:ASN:ND2	2.34	0.55
41:85:66:ASN:HB2	41:85:76:TYR:HB2	1.88	0.55
26:14:929:G:O5'	26:14:929:G:H8	1.90	0.55
36:35:121:LYS:HG3	36:35:122:PRO:HD2	1.88	0.55
26:1H:796:C:H2'	26:1H:797:C:C6	2.41	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:26:G:OP1	43:A5:80:PRO:HB3	2.05	0.55
1:1G:359:U:H2'	1:1G:360:A:C8	2.40	0.55
26:1H:1557:C:OP2	26:1H:1558:A:O2'	2.16	0.55
26:14:2420:C:N4	55:M5:31:HIS:O	2.39	0.55
13:4I:7:VAL:HB	31:4I:115:ARG:CZ	2.36	0.55
1:13:658:G:H2'	1:13:659:U:C6	2.41	0.55
6:5E:97:PHE:HD2	18:9I:31:LEU:HD21	1.72	0.55
29:21:13:ARG:HG2	29:21:13:ARG:HH11	1.72	0.55
30:39:53:THR:HG23	30:39:55:GLY:H	1.71	0.55
1:13:591:U:H2'	1:13:592:G:C8	2.42	0.55
26:1H:1101:U:H2'	26:1H:1102:C:C6	2.42	0.55
22:1K:51:U:H2'	22:1K:52:G:C8	2.40	0.55
26:14:1935:G:H1'	26:14:1964:G:N2	2.22	0.55
26:14:1593:G:H2'	26:14:1594:G:C8	2.41	0.55
29:29:8:LYS:HB3	29:29:192:ASN:HA	1.88	0.55
3:22:150:LYS:HG3	3:22:169:ALA:HB2	1.89	0.55
26:1H:1038:C:H2'	26:1H:1039:G:O4'	2.06	0.55
29:29:49:LEU:O	29:29:78:LEU:HA	2.07	0.55
26:1H:2830:G:H5''	26:1H:2830:G:H8	1.72	0.55
32:51:98:LEU:HD13	32:51:125:VAL:HG23	1.89	0.55
26:1H:1728:G:N1	26:1H:1730:U:OP2	2.38	0.55
29:29:101:ARG:O	29:29:201:THR:OG1	2.25	0.55
27:1J:44:G:H5''	27:1J:45:A:OP1	2.05	0.55
19:AI:41:VAL:HB	19:AI:42:PRO:HA	1.89	0.55
26:14:363(C):G:H2'	26:14:363(D):G:C8	2.42	0.55
26:1H:1533:C:H3'	26:1H:1534:G:C5'	2.35	0.55
1:1G:660:G:H1	1:1G:745:C:H42	1.54	0.55
26:1H:882:G:N2	26:1H:894:C:H42	2.03	0.55
26:1H:569:U:C4	26:1H:570:G:C6	2.95	0.55
1:1G:433:C:H2'	1:1G:434:U:H6	1.71	0.55
26:14:2335:A:C8	26:14:2337:G:C5	2.94	0.55
14:5A:43:CYS:HA	14:5A:46:GLU:OE2	2.07	0.55
34:15:104:LYS:HA	34:15:107:LEU:HD12	1.87	0.55
26:14:1006:C:H1'	34:15:106:MET:HE3	1.89	0.55
26:14:1247:A:OP1	30:39:95:ARG:NH2	2.38	0.55
26:1H:1018:C:O2'	26:1H:1019:U:H5'	2.06	0.55
1:13:1311:G:N2	1:13:1326:C:O2	2.19	0.55
46:H8:73:GLN:HB2	46:H8:87:ASP:HB2	1.87	0.55
1:13:917:G:H2'	1:13:918:A:C8	2.42	0.55
1:1G:316:G:OP2	1:1G:351:G:O2'	2.25	0.55
26:1H:1903:G:OP1	28:11:241:PRO:HB2	2.06	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
11:2A:73:MET:HG2	11:2A:103:LEU:HD13	1.87	0.55
44:F8:36:LYS:HG2	44:F8:54:VAL:HB	1.89	0.55
26:14:751:A:P	58:14:3512:HOH:O	2.62	0.55
26:14:1225:C:O3'	42:95:85:LYS:HD3	2.07	0.55
42:95:85:LYS:CD	42:95:86:GLY:H	2.20	0.55
26:1H:1416:G:HO2'	26:1H:1417:C:H6	1.54	0.55
26:1H:2128:C:N4	26:1H:2160:G:H1	2.03	0.55
55:Q8:14:VAL:HG11	55:Q8:21:LYS:HZ1	1.71	0.55
26:14:362:U:H5'	26:14:363:G:OP2	2.07	0.55
4:32:4:TYR:CE2	4:32:11:LEU:HD11	2.41	0.55
1:1G:464:G:O6	1:1G:466:C:H5'	2.06	0.55
26:1H:1113:U:OP1	32:51:2:SER:N	2.40	0.55
39:A8:84:GLN:HG2	39:A8:110:LEU:HD12	1.89	0.55
26:1H:2636:U:P	29:21:79:ARG:HA	2.46	0.55
30:39:53:THR:HG22	30:39:56:GLU:HG3	1.88	0.55
26:1H:1859:A:N6	26:1H:1883:G:O2'	2.39	0.55
26:1H:1338:G:O2'	26:1H:1339:G:H5'	2.07	0.55
1:1G:1105:A:H2'	1:1G:1106:G:H8	1.71	0.55
1:13:1316:G:N2	1:13:1318:A:H3'	2.20	0.55
46:D5:7:ALA:O	46:D5:8:TYR:CG	2.59	0.55
1:13:137:C:H42	1:13:226:G:H1	1.55	0.55
26:14:424:G:O6	58:14:4028:HOH:O	2.18	0.55
1:1G:77:C:H42	1:1G:92:G:H1	1.53	0.55
26:14:1442:G:H2'	26:14:1443:G:C8	2.42	0.55
26:1H:518:G:H2'	26:1H:519:U:C6	2.42	0.55
4:32:20:TYR:HA	4:32:26:CYS:HB3	1.89	0.55
26:1H:1534:G:H3'	26:1H:1534:G:N3	2.22	0.55
26:14:1040:C:H2'	26:14:1041:C:C6	2.41	0.55
38:55:96:ARG:NH1	38:55:115:GLU:OE1	2.38	0.55
8:72:121:ASP:OD2	8:72:125:ARG:NH2	2.40	0.55
40:75:4:GLY:HA2	40:75:8:LYS:NZ	2.22	0.55
31:41:97:ASP:O	31:41:100:TRP:N	2.40	0.55
26:1H:241:A:H5''	58:1H:4112:HOH:O	2.06	0.55
26:1H:356:G:H2'	26:1H:357:A:H8	1.72	0.55
26:14:330:A:H2	26:14:1210:A:O2'	1.89	0.55
1:13:624:C:H2'	1:13:625:G:H8	1.70	0.55
1:13:1014:A:C2	1:13:1219:U:H1'	2.41	0.55
26:14:734:A:O2'	26:14:1635:G:H5'	2.07	0.55
26:14:2849:U:H4'	26:14:2868:A:C2	2.41	0.55
1:13:1120:G:H2'	1:13:1121:U:C6	2.42	0.55
26:1H:1001:A:H2'	26:1H:1002:G:O4'	2.07	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
15:6I:63:ARG:HG2	15:6I:67:LEU:HD12	1.88	0.55
54:L5:35:ARG:HG3	54:L5:42:LEU:HD11	1.87	0.55
22:1K:37:MIA:H121	22:1K:38:A:N1	2.21	0.55
26:14:141:A:H8	26:14:1595:G:H21	1.53	0.55
7:6E:49:ILE:O	7:6E:53:LYS:HB2	2.07	0.55
26:14:2542:A:H5''	26:14:2542:A:N3	2.22	0.55
26:14:1963:U:H5''	26:14:1963:U:O2	2.07	0.55
26:14:2652:C:H42	26:14:2668:G:H1	1.54	0.55
5:4E:15:ARG:NH1	25:4K:25:A:H3'	2.22	0.55
7:62:111:ARG:NH2	7:62:122:HIS:HB3	2.21	0.55
26:1H:2053:G:H5'	29:21:144:ARG:O	2.07	0.55
26:1H:2503:A:OP1	58:1H:3623:HOH:O	2.18	0.55
45:C5:14:LEU:HD12	45:C5:15:VAL:H	1.71	0.55
1:1G:1208:C:H2'	1:1G:1209:C:H6	1.71	0.55
26:14:399:G:OP2	58:14:3687:HOH:O	2.18	0.55
1:1G:1057:G:H1	1:1G:1203:C:N4	2.03	0.55
45:G8:49:VAL:HG21	45:G8:55:TYR:CE2	2.40	0.55
1:1G:1305:G:H22	1:1G:1331:G:C2'	2.20	0.55
26:1H:270(V):G:H2'	26:1H:270(W):G:C8	2.40	0.55
38:55:77:ARG:HH11	38:55:77:ARG:HG3	1.71	0.55
1:1G:1132:C:H2'	1:1G:1133:G:H8	1.71	0.55
12:3I:8:ASN:O	12:3I:11:VAL:HG23	2.07	0.55
30:39:169:ASN:O	30:39:169:ASN:ND2	2.40	0.55
28:19:16:MET:HE1	28:19:208:LYS:HE2	1.88	0.55
26:14:2062:A:HO2'	26:14:2063:C:P	2.29	0.55
26:14:1027:A:C2	26:14:2488:A:H5'	2.42	0.55
26:14:817:C:H2'	26:14:818:G:O4'	2.07	0.55
1:1G:922:G:H21	1:1G:1398:A:H2	1.55	0.55
1:13:199:G:O6	1:13:218:C:N4	2.40	0.55
26:14:587:C:O2	36:35:33:ARG:NH1	2.40	0.55
42:D8:25:LEU:H	42:D8:92:THR:CG2	2.20	0.55
21:1F:5:ASP:O	21:1F:11:GLY:HA3	2.07	0.55
1:1G:73:G:H1	1:1G:97:U:H3	1.53	0.55
55:Q8:38:GLY:HA2	55:Q8:40:GLU:N	2.22	0.55
53:O8:32:ASN:N	53:O8:32:ASN:OD1	2.39	0.55
26:1H:2341:G:H2'	26:1H:2342:C:C6	2.41	0.55
26:14:1171:G:N2	26:14:1174:A:N1	2.55	0.55
50:L8:13:ILE:HG22	58:L8:201:HOH:O	2.06	0.55
44:F8:1:MET:C	44:F8:3:THR:H	2.10	0.55
26:1H:2057:A:P	58:1H:3613:HOH:O	2.65	0.55
2:12:162:ILE:O	2:12:185:ILE:HG12	2.07	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:13:631:G:C8	1:13:632:A:H2	2.25	0.55
49:K8:50:ILE:HD12	49:K8:51:ARG:H	1.72	0.55
12:3I:58:VAL:O	12:3I:65:GLU:HA	2.07	0.55
41:C8:17:ILE:HG23	41:C8:39:LEU:HD12	1.88	0.55
12:3I:84:LEU:HB2	12:3I:105:TYR:CE2	2.41	0.55
39:A8:14:VAL:O	39:A8:18:ILE:HD13	2.06	0.55
26:14:218:A:H2	26:14:235:U:H4'	1.71	0.55
26:1H:1088:A:H5'	26:1H:1089:G:H5'	1.88	0.55
29:29:39:PRO:HA	29:29:43:GLY:HA2	1.89	0.55
26:14:2693:A:H2'	26:14:2694:G:H8	1.72	0.55
3:22:112:SER:O	3:22:116:VAL:HG23	2.07	0.55
10:1I:48:THR:HA	10:1I:62:HIS:HB3	1.88	0.55
1:13:748:C:O5'	1:13:748:C:H6	1.89	0.55
1:1G:250:A:H1'	1:1G:251:G:OP2	2.07	0.55
27:1J:15:A:H1'	27:1J:109:G:C5	2.42	0.55
51:M8:39:CYS:H	51:M8:41:PRO:CD	2.20	0.55
26:1H:1188:U:O2'	26:1H:1189:A:H5'	2.07	0.55
26:14:39:C:O2	30:39:46:ARG:NH2	2.39	0.55
1:1G:707:C:OP1	11:2A:85:ARG:NH1	2.40	0.55
1:1G:458:C:H2'	1:1G:464:G:H8	1.72	0.55
26:14:731:C:H5''	58:14:3871:HOH:O	2.07	0.55
26:14:2317:C:H2'	26:14:2318:G:O4'	2.06	0.55
33:61:47:LEU:HA	33:61:50:ARG:HB2	1.89	0.55
26:1H:580:C:H2'	26:1H:581:C:H6	1.72	0.55
26:14:729:G:O5'	28:19:208:LYS:NZ	2.38	0.55
27:16:15:A:H1'	27:16:109:G:C8	2.41	0.55
3:2E:128:PHE:HZ	3:2E:132:ARG:HD2	1.71	0.55
50:L8:31:LEU:O	50:L8:32:GLN:HB2	2.06	0.55
26:1H:34:C:H4'	26:1H:35:G:OP1	2.07	0.55
27:1J:21:G:H2'	27:1J:22:U:O4'	2.07	0.55
26:14:1316:U:O2'	26:14:1317:A:H5'	2.07	0.55
33:69:6:LEU:H	33:69:36:ALA:HA	1.72	0.55
26:14:232:G:H8	26:14:232:G:OP2	1.90	0.55
26:1H:516:C:OP1	52:N8:13:LYS:NZ	2.30	0.55
51:M8:18:CYS:SG	51:M8:19:GLY:N	2.78	0.54
1:1G:649:G:H2'	1:1G:650:G:C8	2.43	0.54
36:35:60:MET:C	36:35:61:ARG:HG2	2.27	0.54
11:2A:85:ARG:HA	11:2A:112:THR:OG1	2.07	0.54
26:1H:1798:U:H5'	28:11:259:THR:OG1	2.08	0.54
43:E8:18:ARG:HD3	43:E8:76:VAL:HG13	1.89	0.54
26:14:2439:A:H8	26:14:2439:A:H5'	1.72	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
45:G8:76:CYS:CB	45:G8:97:ARG:HG2	2.37	0.54
26:14:943:U:OP2	36:35:36:LYS:HE3	2.06	0.54
24:3L:5:G:N2	24:3L:68:C:H42	2.04	0.54
29:21:50:GLY:HA2	29:21:77:ILE:HA	1.89	0.54
46:D5:104:PHE:HA	46:D5:139:VAL:H	1.71	0.54
26:1H:2562:U:O2'	35:68:23:ARG:NH1	2.40	0.54
1:1G:181:G:O2'	1:1G:183:G:O6	2.22	0.54
26:1H:1460:A:O2'	26:1H:1461:G:OP1	2.22	0.54
30:39:80:ALA:O	30:39:83:PHE:HB2	2.06	0.54
1:13:237:C:H5''	17:8I:25:ARG:CZ	2.37	0.54
41:C8:58:ARG:HA	41:C8:61:TRP:CE3	2.42	0.54
38:55:44:LEU:HD22	38:55:48:VAL:HG23	1.88	0.54
7:62:65:ALA:HB1	7:62:127:ALA:HB3	1.89	0.54
20:BI:10:LEU:HG	20:BI:12:ALA:H	1.72	0.54
26:14:1111:A:H4'	32:59:3:ARG:HD3	1.89	0.54
42:95:85:LYS:CG	42:95:86:GLY:H	2.21	0.54
36:78:50:ARG:CG	55:Q8:58:ILE:HD11	2.36	0.54
1:1G:909:A:H2'	1:1G:910:C:O4'	2.07	0.54
29:29:151:TYR:HD2	29:29:154:LYS:NZ	2.04	0.54
12:3I:90:VAL:HG12	12:3I:92:ASP:H	1.73	0.54
9:82:112:LYS:HA	9:82:119:ALA:CB	2.37	0.54
31:41:21:ARG:NH1	31:41:21:ARG:HG2	2.20	0.54
1:13:1286:A:C8	1:13:1287:A:H4'	2.43	0.54
52:J5:3:LYS:HD2	52:J5:3:LYS:N	2.21	0.54
35:25:63:VAL:HG12	35:25:106:LEU:HD11	1.89	0.54
26:14:1026:U:H2'	27:1J:88:C:H42	1.72	0.54
2:1E:97:TRP:HZ3	2:1E:172:ILE:HB	1.73	0.54
26:1H:2336:A:H61	47:I8:43:THR:HB	1.71	0.54
26:1H:1167:U:C2	26:1H:1183:G:N2	2.75	0.54
44:F8:65:ARG:HH11	44:F8:70:LEU:HD22	1.71	0.54
26:1H:2855:C:H2'	26:1H:2856:C:H6	1.72	0.54
41:85:88:ILE:HD13	42:95:47:VAL:HG21	1.87	0.54
6:52:24:GLU:OE2	6:52:28:ARG:NH1	2.40	0.54
19:AA:66:MET:N	19:AA:66:MET:SD	2.81	0.54
35:25:49:ARG:HA	35:25:53:LYS:NZ	2.21	0.54
34:58:75:TYR:CE2	34:58:77:GLY:HA2	2.42	0.54
35:25:92:GLU:OE1	35:25:113:LYS:HD2	2.07	0.54
26:1H:322:A:H5'	26:1H:340:A:H1'	1.90	0.54
24:3L:23:A:H2'	24:3L:24:G:H8	1.72	0.54
18:9I:73:ALA:HB1	18:9I:79:LEU:HG	1.89	0.54
1:13:1346:A:H5''	9:8E:120:ARG:NH1	2.22	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
45:G8:87:LYS:O	45:G8:94:LYS:HB2	2.08	0.54
22:1K:9:A:H2	22:1K:11:C:H41	1.54	0.54
1:1G:619:U:N3	4:32:134:ASP:OD1	2.34	0.54
19:AI:41:VAL:O	51:M8:63:TYR:OH	2.26	0.54
30:39:67:GLN:HG3	30:39:67:GLN:O	2.06	0.54
3:22:32:LEU:HD22	3:22:59:ARG:HH12	1.72	0.54
4:3E:107:ARG:NH2	4:3E:194:LEU:HD22	2.23	0.54
7:62:115:ARG:O	7:62:118:VAL:HG22	2.07	0.54
26:14:46:C:H2'	26:14:47:C:C6	2.43	0.54
29:21:60:ASN:OD1	29:21:62:PRO:HD2	2.07	0.54
10:1A:47:PHE:CZ	14:5A:37:PHE:HE2	2.25	0.54
14:5I:21:TYR:HE2	14:5I:23:ARG:NE	2.06	0.54
40:B8:5:ALA:HA	40:B8:8:LYS:HE2	1.90	0.54
26:1H:1636:C:H2'	26:1H:1637:A:C8	2.42	0.54
8:7E:95:VAL:HG12	8:7E:99:GLU:HB2	1.88	0.54
11:2I:86:GLY:N	11:2I:112:THR:OG1	2.27	0.54
8:7E:49:GLU:HG2	8:7E:62:TYR:HE2	1.72	0.54
1:13:156:G:N2	1:13:165:C:N3	2.40	0.54
1:13:963:G:N2	1:13:972:C:N3	2.48	0.54
33:69:69:LYS:NZ	33:69:73:GLU:OE2	2.32	0.54
1:13:1004:A:H2'	1:13:1005:A:O4'	2.06	0.54
1:1G:742:G:P	15:6A:35:ARG:HH22	2.30	0.54
37:88:59:ARG:C	37:88:61:GLY:H	2.09	0.54
26:14:2615:U:C2	52:J5:7:PRO:HA	2.42	0.54
26:1H:2345:G:N3	26:1H:2381:C:H2'	2.22	0.54
40:75:91:ARG:HD3	40:75:120:ARG:HB3	1.89	0.54
5:42:34:VAL:O	5:42:41:VAL:HG12	2.07	0.54
26:14:4:C:H42	26:14:2899:G:N2	2.05	0.54
24:3L:18:G:H2'	24:3L:57:G:N2	2.23	0.54
26:1H:1805:U:O2	28:11:50:THR:HB	2.06	0.54
26:1H:194:G:H2'	26:1H:195:A:O4'	2.08	0.54
26:14:1857:G:O2'	26:14:1885:A:N6	2.40	0.54
26:14:582:G:H2'	26:14:583:G:C8	2.42	0.54
26:1H:818:G:H5'	26:1H:839:U:OP1	2.08	0.54
26:1H:1108:U:O4	26:1H:1109:C:N4	2.41	0.54
1:13:618:C:H5''	1:13:619:U:H5''	1.90	0.54
37:45:17:LEU:HD21	37:45:41:TRP:HE1	1.73	0.54
1:1G:1412:C:H2'	1:1G:1413:A:C8	2.42	0.54
43:E8:60:ASN:OD1	43:E8:60:ASN:N	2.39	0.54
36:78:111:ARG:HG2	36:78:128:HIS:CD2	2.43	0.54
26:1H:1827:C:H2'	26:1H:1828:G:H5'	1.88	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:1H:2593:U:H2'	26:1H:2594:C:H6	1.71	0.54
29:29:68:ALA:C	29:29:70:ALA:H	2.10	0.54
55:M5:48:PHE:CE2	55:M5:50:LEU:HD13	2.42	0.54
13:4A:37:THR:HG22	13:4A:55:ARG:HD2	1.90	0.54
1:13:1133:G:H22	1:13:1141:C:N4	2.04	0.54
30:39:7:TYR:CE2	30:39:10:PRO:HG3	2.42	0.54
39:A8:24:LEU:HB2	39:A8:85:VAL:HG12	1.88	0.54
26:1H:1069:A:O2'	26:1H:1072:C:OP1	2.22	0.54
29:29:202:LYS:N	29:29:202:LYS:HD2	2.22	0.54
26:14:527:C:OP2	26:14:2779:U:H5	1.90	0.54
1:1G:1130:A:N6	1:1G:1144:G:H21	2.05	0.54
55:Q8:34:TRP:CD1	55:Q8:35:GLN:N	2.75	0.54
26:1H:994:C:OP1	41:C8:53:ARG:NH2	2.41	0.54
39:A8:99:LYS:HE2	39:A8:103:GLU:OE1	2.07	0.54
55:Q8:38:GLY:HA2	55:Q8:41:ILE:H	1.73	0.54
26:14:46:C:H2'	26:14:47:C:H6	1.72	0.54
26:1H:1058:U:H3	26:1H:1080:A:H61	1.54	0.54
26:14:2228:G:OP1	28:19:261:LYS:NZ	2.38	0.54
24:3L:58:A:H2'	24:3L:60:U:OP2	2.07	0.54
26:14:2257:U:H2'	26:14:2258:C:C6	2.42	0.54
26:1H:768:G:N7	58:1H:4260:HOH:O	2.33	0.54
26:1H:1221:C:H2'	26:1H:1222:C:H6	1.72	0.54
5:4E:147:ASP:HA	5:4E:150:ARG:HH22	1.71	0.54
36:35:15:ARG:CZ	36:35:15:ARG:HB2	2.37	0.54
51:M8:18:CYS:SG	51:M8:39:CYS:HB3	2.47	0.54
26:14:1111:A:O2'	32:59:2:SER:OG	2.25	0.54
26:14:1485:G:H2'	26:14:1486:A:C8	2.43	0.54
27:1J:66:A:N6	27:1J:107:U:H2'	2.23	0.54
40:B8:26:ASP:HB3	40:B8:92:GLY:H	1.72	0.54
30:39:25:PRO:HG3	30:39:28:ILE:HG23	1.89	0.54
26:14:958:U:O2	27:1J:89(A):A:O2'	2.23	0.54
1:13:976:G:H5'	1:13:1358:U:O2'	2.07	0.54
26:1H:962:G:H2'	26:1H:963:U:C6	2.43	0.54
33:61:131:LYS:HB3	33:61:132:PRO:HA	1.89	0.54
26:1H:507:A:H5''	26:1H:508:G:H3'	1.89	0.54
41:85:65:ILE:HD11	41:85:96:ALA:HB3	1.88	0.54
9:8E:27:THR:N	9:8E:61:ALA:O	2.36	0.54
26:1H:2887:U:H2'	26:1H:2888:C:H6	1.72	0.54
35:25:63:VAL:HB	35:25:102:VAL:HG12	1.90	0.54
10:1A:4:ILE:HG12	10:1A:100:THR:HG22	1.89	0.54
1:13:17:U:H2'	1:13:18:C:C6	2.42	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
41:85:16:LYS:O	41:85:20:LEU:HD23	2.08	0.54
46:D5:118:GLN:HG3	46:D5:173:ALA:O	2.07	0.54
1:1G:542:G:H5'	4:32:41:GLY:HA3	1.88	0.54
5:4E:11:ILE:HD11	5:4E:31:LEU:HD13	1.90	0.54
46:D5:97:GLU:HB3	46:D5:125:LEU:HD11	1.90	0.54
52:J5:41:PRO:O	52:J5:44:THR:OG1	2.24	0.54
46:D5:110:GLY:HA2	46:D5:144:LEU:H	1.72	0.54
1:1G:688:G:H2'	1:1G:689:C:H6	1.73	0.54
1:1G:160:A:H1'	1:1G:344:A:C5	2.43	0.54
45:G8:9:LYS:HA	45:G8:27:VAL:HG22	1.88	0.54
55:Q8:53:PRO:HB2	55:Q8:56:GLU:HB2	1.89	0.54
26:14:251:A:C5	26:14:252:G:H1'	2.42	0.54
26:14:1771:C:HO2'	26:14:1786:A:H8	1.54	0.54
1:13:875:C:O2'	8:7E:14:ARG:NH1	2.41	0.54
9:8E:45:ALA:HA	9:8E:48:GLU:HG2	1.89	0.54
26:1H:2116:G:O6	26:1H:2172:U:N3	2.40	0.54
39:65:66:ALA:O	39:65:70:GLY:N	2.33	0.54
14:5I:3:ARG:O	14:5I:7:ILE:HG23	2.08	0.54
26:1H:1077:A:H3'	26:1H:1078:U:C5'	2.38	0.54
1:13:1095:U:H5'	1:13:1109:C:O2	2.08	0.54
47:E5:12:ASN:HA	47:E5:14:ARG:HH21	1.72	0.54
1:1G:1375:A:H4'	7:62:29:LYS:HE3	1.88	0.54
34:58:42:TRP:HA	34:58:48:MET:CE	2.37	0.54
32:59:9:ILE:HG22	32:59:51:ARG:HA	1.88	0.54
26:14:2789:C:H2'	26:14:2790:A:O4'	2.08	0.54
19:AA:63:THR:OG1	19:AA:65:ASN:O	2.21	0.54
28:11:89:SER:HB2	28:11:159:ALA:CB	2.37	0.54
3:2E:150:LYS:HB3	3:2E:201:TYR:HB2	1.89	0.54
53:K5:36:LEU:HD23	53:K5:50:ARG:HD3	1.89	0.54
54:P8:12:ARG:HH21	54:P8:44:PRO:HB3	1.73	0.54
36:35:99:LEU:HD12	36:35:102:ARG:HH22	1.73	0.54
26:1H:1593:G:H2'	26:1H:1594:G:C8	2.43	0.54
26:14:2105:C:H42	26:14:2184:G:H1	1.55	0.54
27:1J:116:G:H8	27:1J:116:G:O5'	1.90	0.54
1:1G:944:G:O2'	1:1G:1339:A:N6	2.40	0.54
4:3E:152:SER:HB2	4:3E:155:LEU:HG	1.90	0.54
31:41:10:LYS:O	31:41:15:VAL:HG23	2.07	0.54
17:8A:56:VAL:HG12	17:8A:77:VAL:HB	1.90	0.54
2:1E:18:GLY:N	2:1E:42:ILE:HG22	2.22	0.54
26:1H:899:A:HO2'	26:1H:900:A:H8	1.54	0.54
24:3K:18:G:O2'	24:3K:19:G:OP1	2.24	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:800:A:P	58:14:3859:HOH:O	2.64	0.54
1:13:1157:A:N6	1:13:1178:G:H21	2.06	0.54
26:1H:2377:A:H4'	39:A8:111:GLU:HB3	1.90	0.54
1:1G:1210:C:H3'	1:1G:1211:U:H5''	1.88	0.54
1:13:1098:C:C2	1:13:1099:G:C8	2.96	0.54
39:65:84:GLN:HB3	39:65:110:LEU:H	1.73	0.54
1:13:277:C:P	17:8I:68:ARG:HH12	2.30	0.54
26:14:2776:A:OP1	26:14:2776:A:H3'	2.07	0.54
20:BI:49:ALA:O	20:BI:52:ALA:N	2.40	0.54
12:3A:111:LYS:HD2	12:3A:111:LYS:H	1.72	0.54
27:16:31:C:H2'	27:16:32:C:C6	2.41	0.54
28:11:149:PRO:O	28:11:150:LYS:HB2	2.08	0.54
43:E8:45:TYR:CZ	43:E8:49:LYS:HE3	2.43	0.54
53:O8:47:THR:HG22	53:O8:48:VAL:H	1.72	0.54
1:1G:1244:C:H2'	1:1G:1245:A:C8	2.43	0.54
43:E8:27:LYS:HB3	43:E8:31:GLU:HG3	1.90	0.54
32:51:77:LYS:HE2	32:51:138:LYS:HB2	1.90	0.54
31:49:160:VAL:HG12	31:49:161:THR:H	1.73	0.54
3:2E:3:ASN:OD1	3:2E:3:ASN:N	2.40	0.54
5:42:93:PRO:HG2	8:72:105:ARG:NE	2.22	0.54
1:1G:713:G:H2'	1:1G:714:G:C8	2.42	0.54
2:1E:163:PHE:CD1	2:1E:185:ILE:HG13	2.43	0.54
32:59:6:ARG:HB2	32:59:66:GLY:HA2	1.90	0.54
27:16:42:C:O2'	31:41:67:LYS:HE3	2.08	0.54
45:C5:47:LYS:HG3	45:C5:60:PHE:CE2	2.43	0.54
1:1G:502:G:OP1	12:3A:118:SER:HB3	2.08	0.54
1:1G:838:G:H2'	1:1G:841:U:H5''	1.89	0.54
1:1G:1371:G:OP1	9:82:11:LYS:HB3	2.08	0.54
52:J5:46:CYS:SG	52:J5:48:GLU:HG2	2.48	0.54
4:3E:148:VAL:HG12	4:3E:149:ALA:O	2.08	0.54
26:1H:2830:G:H5''	26:1H:2830:G:C8	2.43	0.54
10:1I:6:ILE:HG22	10:1I:98:ILE:HG13	1.89	0.54
22:1K:5:G:H1'	22:1K:69:G:N2	2.23	0.54
1:13:1422:G:H5'	35:68:48:PRO:HB3	1.89	0.54
41:C8:44:ASN:ND2	42:D8:75:PHE:O	2.40	0.54
1:13:677:U:H3	1:13:713:G:H22	1.55	0.54
43:A5:62:HIS:HB2	43:A5:64:MET:HG3	1.89	0.54
26:1H:1970:A:P	58:1H:4211:HOH:O	2.62	0.54
26:1H:1359:A:H2	26:1H:1372:U:O4	1.91	0.54
1:13:536:C:H2'	1:13:537:G:C8	2.43	0.54
26:1H:1510:A:N3	26:1H:1510:A:H2'	2.22	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:78:59:LEU:HD11	55:Q8:10:ALA:HA	1.89	0.54
44:B5:8:ILE:HD12	44:B5:8:ILE:H	1.72	0.54
1:13:963:G:H21	10:1I:55:LYS:NZ	2.06	0.54
26:14:2392:A:C8	36:35:61:ARG:HD2	2.43	0.54
46:D5:157:LEU:CB	46:D5:161:VAL:HG11	2.35	0.54
50:L8:50:VAL:HG23	50:L8:54:VAL:HG11	1.89	0.54
26:1H:1494:A:H2'	26:1H:1495:A:C8	2.43	0.54
1:1G:1249:C:O2'	9:82:73:GLN:OE1	2.25	0.54
1:13:353:A:C8	1:13:353:A:H5'	2.38	0.54
26:1H:2680:C:OP2	29:21:111:ARG:NH2	2.41	0.54
1:1G:1367:C:H5'	10:1A:60:ARG:NH2	2.23	0.54
2:1E:15:VAL:HB	2:1E:210:SER:HB2	1.89	0.54
49:K8:48:HIS:N	49:K8:50:ILE:HD11	2.23	0.54
1:1G:942:G:H2'	1:1G:943:U:H6	1.72	0.54
26:1H:2327:A:H2'	26:1H:2328:A:H8	1.72	0.54
29:29:188:VAL:HG23	29:29:189:PRO:HD2	1.89	0.54
1:1G:1224:G:O2'	1:1G:1322:C:OP2	2.26	0.54
33:61:46:ALA:C	33:61:50:ARG:HD3	2.28	0.54
16:7I:11:SER:HB2	16:7I:14:ASN:HB3	1.89	0.54
26:1H:198:C:O2'	26:1H:199:A:H5'	2.08	0.54
46:D5:139:VAL:HG22	46:D5:156:LYS:HG2	1.89	0.54
5:42:67:VAL:HG21	5:42:140:ARG:HA	1.90	0.54
27:1J:87:G:H3'	27:1J:88:C:C5'	2.38	0.54
34:58:133:GLN:O	34:58:134:ARG:HB2	2.07	0.54
26:1H:2146:C:H4'	26:1H:2147:G:N7	2.23	0.54
6:52:23:LYS:HE3	6:52:61:LEU:HD21	1.90	0.54
44:B5:65:ARG:HB3	44:B5:70:LEU:HB3	1.89	0.54
26:1H:1971:A:C4	28:11:241:PRO:HD3	2.43	0.54
19:AA:66:MET:N	19:AA:67:VAL:HB	2.23	0.54
26:1H:2108:C:H2'	26:1H:2109:U:O4'	2.07	0.54
26:1H:1058:U:H2'	26:1H:1059:G:H8	1.73	0.54
1:13:1190:G:OP1	3:2E:4:LYS:HA	2.08	0.54
1:13:474:G:H5''	16:7I:81:ARG:HE	1.72	0.54
26:14:1190:G:H2'	26:14:1191:G:H8	1.73	0.54
26:1H:638:G:C5	26:1H:651:G:C2	2.96	0.54
1:13:760:G:H2'	1:13:761:G:H5'	1.89	0.54
6:52:25:ILE:HD13	6:52:82:ARG:HD2	1.90	0.54
33:61:1:MET:O	33:61:20:ASP:HA	2.07	0.54
32:59:58:GLU:HB2	32:59:61:HIS:ND1	2.23	0.54
26:14:1999:C:H4'	26:14:2723:C:O2	2.07	0.54
26:1H:832:G:H5'	36:78:45:LEU:HD11	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:68:98:VAL:HG11	35:68:114:ILE:HG23	1.89	0.54
1:13:135:C:H2'	1:13:136:C:H5'	1.89	0.54
26:1H:1026:U:H1'	26:1H:1027:A:O5'	2.07	0.54
1:1G:719:C:O2'	18:9A:49:LYS:HB3	2.08	0.54
5:4E:65:ASN:OD1	5:4E:140:ARG:NH2	2.41	0.54
43:E8:68:ARG:O	43:E8:110:LYS:HB2	2.07	0.54
26:1H:2137:C:H42	26:1H:2154:G:N2	2.05	0.53
48:J8:92:LYS:HA	48:J8:95:LEU:HB2	1.89	0.53
1:13:1150:U:O2	10:1I:39:PRO:HG2	2.07	0.53
7:6E:16:LEU:HD12	9:8E:42:ARG:HA	1.90	0.53
1:13:142:G:H2'	1:13:143:A:C8	2.43	0.53
16:7I:39:TYR:CZ	16:7I:41:PRO:HB3	2.44	0.53
26:1H:825:C:O2	36:78:55:ARG:HD3	2.09	0.53
2:12:77:ALA:O	2:12:81:VAL:HG23	2.07	0.53
1:1G:448:A:P	1:1G:485:G:H22	2.31	0.53
2:1E:7:VAL:HG21	2:1E:217:ARG:HH11	1.72	0.53
1:13:109:A:N7	1:13:326:G:H2'	2.23	0.53
1:13:688:G:H2'	1:13:689:C:H6	1.72	0.53
47:I8:23:VAL:HA	47:I8:38:VAL:HG22	1.90	0.53
35:25:4:PRO:O	35:25:5:GLN:HB2	2.08	0.53
3:22:156:ARG:NH2	3:22:159:GLY:O	2.41	0.53
1:1G:1123:A:O2'	10:1A:37:PRO:O	2.18	0.53
5:4E:51:VAL:O	5:4E:55:VAL:HG23	2.08	0.53
3:2E:24:ALA:HB1	3:2E:28:GLN:HB2	1.90	0.53
26:14:1417:C:H42	26:14:1581:G:H1	1.56	0.53
45:G8:84:ARG:O	45:G8:84:ARG:HD2	2.08	0.53
26:1H:1508:A:O2'	26:1H:1509:C:O4'	2.26	0.53
26:14:1997:G:H5''	58:14:3782:HOH:O	2.06	0.53
1:1G:456:C:N4	1:1G:476:G:H1	2.04	0.53
31:41:64:THR:HG23	31:41:94:LEU:HD22	1.91	0.53
26:14:2572:A:N7	29:29:145:LYS:HB2	2.23	0.53
26:14:2572:A:H62	29:29:145:LYS:HD2	1.72	0.53
26:1H:1061:U:O2'	26:1H:1070:A:O4'	2.25	0.53
1:1G:1028(A):C:H5	1:1G:1029:G:C5	2.25	0.53
26:14:249:C:O2	55:M5:12:LYS:NZ	2.38	0.53
23:2L:26:C:H2'	23:2L:27:G:O4'	2.09	0.53
27:1J:90:C:H5'	37:45:18:LYS:HA	1.90	0.53
9:8E:125:TYR:HD1	9:8E:126:SER:N	2.06	0.53
26:1H:1889:A:N1	26:1H:2234:G:H1'	2.23	0.53
1:13:474:G:H2'	1:13:475:G:C8	2.43	0.53
1:1G:9:G:OP1	5:42:122:GLU:HB2	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
27:1J:94:C:H2'	27:1J:95:U:C6	2.44	0.53
31:41:139:LEU:HD22	31:41:146:TYR:HD1	1.73	0.53
6:52:7:ASN:N	6:52:7:ASN:OD1	2.42	0.53
19:AI:6:LYS:O	19:AI:7:LYS:HB3	2.07	0.53
7:62:95:ARG:HH21	7:62:99:LEU:HD11	1.74	0.53
28:19:239:ARG:NE	58:19:302:HOH:O	2.41	0.53
1:13:67:C:H2'	1:13:68:G:C8	2.44	0.53
26:1H:646:A:H2'	26:1H:647:G:O4'	2.09	0.53
27:16:54:G:H2'	27:16:55:U:H6	1.73	0.53
26:1H:1901:A:OP2	28:11:255:LYS:HE2	2.08	0.53
15:6A:4:THR:O	15:6A:7:GLU:HG2	2.07	0.53
17:8A:63:ARG:HG2	17:8A:64:PRO:HD2	1.89	0.53
26:1H:2074:U:H2'	26:1H:2075:U:C6	2.44	0.53
26:1H:1871:A:H2'	26:1H:1872:A:C8	2.42	0.53
26:1H:249:C:P	58:1H:3665:HOH:O	2.64	0.53
1:1G:1502:A:H5''	1:1G:1504:G:C8	2.43	0.53
26:14:2748:A:H2'	26:14:2749:A:C8	2.43	0.53
46:D5:4:ARG:HG2	46:D5:58:VAL:HG21	1.88	0.53
26:1H:1534:G:H22	26:1H:1538:G:H1	1.56	0.53
26:14:2376:A:C2	39:65:112:PHE:HB2	2.43	0.53
28:11:145:VAL:HG12	28:11:146:GLU:O	2.08	0.53
26:14:690:G:O2'	28:19:43:ARG:NH2	2.41	0.53
26:14:1538:G:H2'	26:14:1539:G:H8	1.74	0.53
33:69:79:ILE:HD11	33:69:140:LEU:HD11	1.91	0.53
55:Q8:33:ASN:O	55:Q8:34:TRP:CD1	2.62	0.53
26:1H:363:G:H2'	26:1H:363(A):A:H8	1.73	0.53
26:1H:270(P):C:H2'	26:1H:270(Q):C:C6	2.42	0.53
1:1G:1478:C:H2'	1:1G:1479:C:H6	1.73	0.53
26:14:2052:G:O4'	29:29:142:GLY:HA3	2.08	0.53
7:62:18:TYR:HB3	7:62:59:LEU:HD12	1.90	0.53
26:1H:1932:A:H2'	26:1H:1933:G:O4'	2.08	0.53
33:61:107:VAL:HG12	33:61:108:THR:H	1.73	0.53
41:85:85:LYS:HD3	41:85:116:ALA:O	2.09	0.53
31:41:33:ARG:O	31:41:162:THR:HG23	2.08	0.53
26:1H:1728:G:H3'	26:1H:1729:A:H5''	1.88	0.53
55:M5:34:TRP:CD1	55:M5:35:GLN:N	2.74	0.53
26:14:2873:A:H8	38:55:6:SER:N	2.06	0.53
1:13:1423:G:P	35:68:49:ARG:HH22	2.31	0.53
26:1H:2318:G:H22	39:A8:2:ALA:N	2.07	0.53
26:1H:1453:A:O2'	26:1H:1454:U:H2'	2.07	0.53
40:75:24:PRO:HA	40:75:49:VAL:HG23	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
29:21:34:VAL:HG22	29:21:48:GLN:HB3	1.90	0.53
1:13:262:A:H5''	1:13:263:A:OP2	2.08	0.53
32:51:8:PRO:O	32:51:10:PRO:HD3	2.07	0.53
13:4A:60:VAL:HG13	13:4A:64:TRP:HE1	1.73	0.53
15:6I:39:LEU:HD13	15:6I:56:LEU:HB2	1.90	0.53
26:14:582:G:N7	58:14:3964:HOH:O	2.34	0.53
26:1H:2675:A:H5'	35:68:29:ASN:O	2.09	0.53
1:1G:284:G:H2'	1:1G:285:G:H8	1.73	0.53
44:B5:55:ASN:HB2	44:B5:80:ILE:HG13	1.89	0.53
33:69:102:SER:HB2	33:69:108:THR:HG23	1.90	0.53
1:1G:952:U:H4'	1:1G:964:A:N1	2.23	0.53
44:F8:15:GLU:HG3	44:F8:16:LYS:N	2.23	0.53
33:61:55:ALA:HA	33:61:58:LEU:HB3	1.90	0.53
26:1H:2023:G:H5'	26:1H:2617:C:H4'	1.89	0.53
7:62:69:VAL:HG12	7:62:103:TRP:HE3	1.73	0.53
1:1G:789:U:H2'	1:1G:791:G:N7	2.24	0.53
45:G8:29:GLU:HB3	45:G8:38:ILE:CG2	2.38	0.53
33:61:81:VAL:HG11	33:61:88:ILE:HD12	1.91	0.53
26:1H:2032:G:N2	29:21:146:THR:HG23	2.20	0.53
1:13:10:A:OP2	5:4E:126:ARG:HD3	2.08	0.53
8:7E:83:ILE:HB	8:7E:137:VAL:HG13	1.91	0.53
10:1I:57:LYS:HD2	10:1I:60:ARG:NH1	2.22	0.53
16:7I:53:VAL:O	16:7I:57:ARG:HG2	2.08	0.53
13:4A:14:ARG:HD2	13:4A:42:ALA:HA	1.91	0.53
28:19:255:LYS:H	28:19:255:LYS:NZ	2.06	0.53
46:H8:146:ILE:HD13	46:H8:174:VAL:HG12	1.90	0.53
29:21:78:LEU:O	29:21:79:ARG:HB2	2.08	0.53
2:1E:5:ILE:HG13	2:1E:6:THR:N	2.24	0.53
28:11:182:LEU:H	28:11:272:ALA:CB	2.22	0.53
26:1H:1055:G:H1'	26:1H:1085:A:C2	2.44	0.53
26:14:1161:C:H2'	26:14:1162:G:C8	2.44	0.53
1:13:22:G:H2'	1:13:23:C:C6	2.44	0.53
18:9I:66:LEU:O	18:9I:70:ILE:HG13	2.07	0.53
6:52:7:ASN:HD22	18:9A:76:LEU:HD11	1.74	0.53
6:5E:86:ARG:O	6:5E:87:ARG:HG2	2.08	0.53
1:1G:1059:C:OP2	3:22:199:LYS:NZ	2.41	0.53
1:13:393:A:OP2	16:7I:12:LYS:NZ	2.30	0.53
47:I8:29:GLN:H	47:I8:67:VAL:HG23	1.74	0.53
55:M5:29:LYS:HB3	55:M5:44:LYS:CB	2.38	0.53
40:75:124:ASP:O	40:75:128:GLU:HB2	2.08	0.53
3:2E:62:ASP:O	3:2E:98:ASN:ND2	2.42	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
30:31:152:GLU:HB2	30:31:191:ARG:HD2	1.91	0.53
37:88:66:ILE:HG22	37:88:67:ARG:N	2.24	0.53
26:14:443:A:H1'	26:14:1201:C:O4'	2.09	0.53
8:72:51:VAL:HG11	8:72:60:ARG:HE	1.74	0.53
1:13:554:C:H2'	1:13:555:C:H6	1.72	0.53
17:81:88:TYR:HD2	17:81:89:LEU:HD22	1.73	0.53
40:75:20:PRO:HD2	40:75:86:ILE:HG23	1.89	0.53
30:39:117:ARG:HH12	36:35:1:MET:N	2.04	0.53
1:13:974:A:OP2	14:51:29:ARG:NH2	2.40	0.53
19:AI:41:VAL:HG11	19:AI:67:VAL:HA	1.90	0.53
45:C5:76:CYS:SG	45:C5:97:ARG:HG3	2.49	0.53
26:1H:1257:C:H4'	30:31:83:PHE:CE1	2.43	0.53
49:K8:55:ARG:O	49:K8:58:ALA:HB3	2.08	0.53
26:14:882:G:H22	26:14:894:C:N4	2.07	0.53
1:1G:426:G:OP1	4:32:38:TYR:OH	2.22	0.53
26:14:1533:C:C4	26:14:1534:G:H1'	2.44	0.53
26:1H:831:G:N7	58:1H:3980:HOH:O	2.33	0.53
1:13:590:C:N3	1:13:649:G:N2	2.52	0.53
1:13:881:G:P	12:31:12:ARG:HH22	2.31	0.53
26:14:445:C:O2'	26:14:446:G:H5'	2.08	0.53
26:1H:547:A:H8	26:1H:547:A:O5'	1.92	0.53
15:6I:39:LEU:HB3	15:6I:56:LEU:HD12	1.89	0.53
1:1G:600:C:H2'	1:1G:601:C:H6	1.73	0.53
36:35:122:PRO:HB3	36:35:141:ALA:HB1	1.90	0.53
36:78:39:LYS:HG3	36:78:45:LEU:HD22	1.89	0.53
26:14:192:C:P	58:14:3625:HOH:O	2.66	0.53
29:29:116:VAL:HG23	29:29:120:TRP:HD1	1.74	0.53
35:25:47:ILE:HG23	35:25:48:PRO:HD2	1.91	0.53
26:14:952:G:C6	26:14:966:G:C6	2.96	0.53
26:14:676:A:H8	26:14:2069:G:N2	1.90	0.53
1:1G:649:G:H2'	1:1G:650:G:H8	1.72	0.53
33:69:81:VAL:O	33:69:143:SER:OG	2.26	0.53
5:42:43:LEU:HD22	5:42:136:MET:HG3	1.90	0.53
5:42:136:MET:O	5:42:140:ARG:N	2.32	0.53
26:1H:2689:U:H5''	26:1H:2713:A:H2	1.73	0.53
26:1H:2144:U:N3	26:1H:2146:C:O2	2.42	0.53
15:6A:56:LEU:O	15:6A:60:VAL:HG23	2.09	0.53
26:1H:2788:C:O2'	26:1H:2809:A:N3	2.40	0.53
26:14:2677:G:H2'	26:14:2678:C:C6	2.43	0.53
33:61:57:ARG:O	33:61:61:ARG:HG2	2.09	0.53
1:1G:811:C:O2'	1:1G:901:A:N1	2.41	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
39:A8:25:ARG:NH1	39:A8:42:ASP:OD2	2.41	0.53
28:19:141:VAL:HG23	28:19:162:SER:HB2	1.90	0.53
6:52:68:PRO:HG2	6:52:71:ARG:HG3	1.90	0.53
3:2E:134:ILE:HG22	3:2E:168:ALA:HB3	1.90	0.53
26:1H:232:G:OP2	26:1H:232:G:H8	1.91	0.53
36:78:126:VAL:HG12	36:78:147:LEU:HD22	1.90	0.53
26:14:1899:G:N2	26:14:1902:C:N4	2.49	0.53
26:14:832:G:H5'	36:35:45:LEU:CD1	2.38	0.53
42:95:87:HIS:CE1	42:95:89:GLN:HB2	2.44	0.53
26:1H:1021:A:H8	26:1H:1022:G:H5''	1.74	0.53
24:3L:64:A:C2	24:3L:65:G:H1'	2.44	0.53
26:1H:1049:C:H2'	26:1H:1050:A:H5'	1.90	0.53
5:4E:100:VAL:HG13	5:4E:118:ILE:HG22	1.89	0.53
13:4I:12:ASN:OD1	13:4I:13:LYS:N	2.41	0.53
17:8I:29:HIS:CE1	17:8I:32:TYR:HD2	2.26	0.53
26:1H:910:A:C5	37:88:13:GLN:HG3	2.44	0.53
37:45:75:THR:HB	37:45:86:GLY:HA3	1.91	0.53
55:Q8:32:LEU:HG	55:Q8:33:ASN:N	2.24	0.53
20:BI:10:LEU:HD11	20:BI:12:ALA:HB3	1.89	0.53
3:2E:28:GLN:O	3:2E:31:HIS:N	2.41	0.53
1:13:107:G:O6	20:BI:15:ARG:HG3	2.09	0.53
53:K5:14:THR:OG1	53:K5:19:ARG:HA	2.08	0.53
7:6E:94:ARG:O	7:6E:97:GLN:HB3	2.09	0.53
1:13:1147:C:O2	9:8E:16:ARG:NH1	2.42	0.53
2:1E:136:VAL:O	2:1E:140:HIS:N	2.42	0.53
3:22:90:GLU:HA	3:22:93:LYS:HB2	1.91	0.53
27:16:78:A:C2	27:16:99:A:C4	2.96	0.53
48:F5:41:ARG:HD3	48:F5:43:TYR:HE1	1.73	0.53
22:1K:47:U:H5'	22:1K:48:C:H5'	1.90	0.53
1:13:1106:G:H5''	3:2E:172:ARG:HG2	1.90	0.53
10:1A:30:SER:OG	10:1A:81:THR:HG22	2.08	0.53
1:13:49:U:C2	1:13:361:G:N2	2.76	0.53
26:14:265:A:H1'	26:14:266:G:O4'	2.09	0.53
29:29:56:PRO:HD2	29:29:58:ARG:NH2	2.24	0.53
23:2L:22:A:N6	23:2L:47:7MG:H2'	2.22	0.53
26:1H:588:U:H2'	26:1H:589:C:C6	2.44	0.53
27:1J:13:A:H5''	27:1J:15:A:C6	2.43	0.53
2:1E:163:PHE:HD1	2:1E:185:ILE:HG13	1.73	0.53
1:13:926:G:N2	25:4K:15:A:OP2	2.36	0.53
36:78:59:LEU:O	55:Q8:13:ARG:HD2	2.08	0.53
26:1H:1416:G:H1	26:1H:1582:C:H42	1.56	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
31:49:42:GLY:O	31:49:43:LEU:HD13	2.09	0.53
4:32:24:GLU:OE2	4:32:24:GLU:N	2.41	0.53
1:13:1157:A:H1'	1:13:1158:C:C4	2.44	0.53
2:12:73:THR:HB	2:12:96:ARG:H	1.74	0.53
19:AI:40:ILE:HD11	19:AI:62:ILE:HD13	1.89	0.53
26:14:2308:G:H3'	26:14:2310:A:OP2	2.08	0.53
37:45:25:ASP:CB	37:45:102:VAL:H	2.22	0.53
1:1G:371:G:H1	1:1G:390:C:H42	1.56	0.53
26:14:140:A:C8	26:14:1408:C:O2'	2.62	0.53
3:22:37:GLN:O	3:22:40:ARG:HG3	2.09	0.53
26:14:2027:G:H2'	26:14:2028:U:O4'	2.09	0.53
1:13:1326:C:H2'	1:13:1327:C:C6	2.44	0.53
1:13:1327:C:OP2	21:1F:12:LYS:NZ	2.40	0.53
50:L8:26:LEU:HB2	50:L8:28:LEU:HD12	1.89	0.53
50:L8:4:LEU:HG	50:L8:39:ASP:HB2	1.91	0.53
26:14:460:A:H5''	26:14:461:C:OP2	2.09	0.53
19:AA:42:PRO:HG3	51:I5:60:GLN:HG3	1.91	0.53
1:13:1080:A:H5''	5:4E:16:THR:HG21	1.91	0.53
1:13:1077:G:N2	1:13:1080:A:OP2	2.31	0.53
42:95:21:ARG:NH2	42:95:91:TYR:O	2.41	0.53
1:1G:1307:U:H2'	1:1G:1308:U:C6	2.44	0.53
28:19:182:LEU:H	28:19:272:ALA:HB2	1.73	0.53
17:8I:45:HIS:O	17:8I:73:VAL:HG23	2.09	0.53
26:1H:309:G:N3	26:1H:329:G:O2'	2.41	0.53
17:8I:58:GLU:O	17:8I:74:LEU:N	2.35	0.53
26:1H:2068:U:H3	26:1H:2430:A:H2	1.55	0.53
26:1H:150:C:H2'	26:1H:151:C:C6	2.44	0.53
1:1G:1281:U:H3'	1:1G:1282:C:C5	2.44	0.53
26:1H:1359:A:H2'	26:1H:1360:A:H5'	1.91	0.53
26:1H:2061:G:OP2	26:1H:2502:G:H5'	2.09	0.53
26:1H:2502:G:P	58:1H:3635:HOH:O	2.66	0.53
26:14:2270:G:OP2	58:14:3726:HOH:O	2.19	0.53
26:14:782:A:H5'	26:14:783:A:C2	2.43	0.53
26:14:2685:G:OP2	40:75:51:ARG:NH1	2.42	0.53
26:14:93:C:H4'	45:C5:54:LYS:HE3	1.90	0.53
40:75:1:MET:HB3	40:75:6:LEU:HB2	1.90	0.53
29:29:11:MET:HE3	29:29:186:GLY:HA2	1.91	0.53
4:3E:102:ASP:HB3	4:3E:136:PRO:CB	2.39	0.53
13:4I:82:MET:O	13:4I:84:ILE:N	2.42	0.53
1:13:260:G:H2'	1:13:261:U:C6	2.44	0.53
12:3A:84:LEU:HD12	12:3A:104:VAL:HG11	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
54:P8:15:THR:HG22	54:P8:16:HIS:CE1	2.43	0.53
39:A8:70:GLY:O	39:A8:105:ALA:HA	2.09	0.53
4:32:106:TYR:HE1	4:32:112:VAL:O	1.91	0.53
26:14:2303:G:N2	26:14:2314:C:N3	2.56	0.53
1:1G:1082:G:OP2	1:1G:1082:G:H8	1.91	0.53
26:1H:931:G:O2'	50:L8:24:LYS:NZ	2.35	0.53
51:I5:21:VAL:HG22	51:I5:22:ILE:H	1.74	0.52
26:1H:674:G:H1'	30:31:74:ARG:HD3	1.90	0.52
31:49:43:LEU:HD12	31:49:45:GLU:OE1	2.08	0.52
9:8E:97:LYS:HB2	9:8E:102:LEU:HD12	1.90	0.52
4:32:18:LYS:HD2	4:32:20:TYR:CZ	2.44	0.52
41:85:28:ARG:HH11	41:85:38:THR:HG1	1.54	0.52
33:61:41:GLU:O	33:61:45:LYS:HB2	2.09	0.52
29:21:105:THR:HB	29:21:197:ILE:HG23	1.91	0.52
32:51:4:ILE:HD11	32:51:7:LEU:HD11	1.90	0.52
2:12:178:ARG:HH21	8:72:68:ARG:NH2	2.07	0.52
2:1E:178:ARG:HB2	2:1E:178:ARG:HH11	1.74	0.52
1:1G:1261:A:H5'	1:1G:1283:G:O3'	2.09	0.52
9:8E:18:PHE:HD2	9:8E:62:TYR:HD2	1.57	0.52
26:1H:1230:C:H2'	26:1H:1231:G:H8	1.71	0.52
1:1G:1080:A:H5'	5:42:14:ARG:HH22	1.73	0.52
26:14:1257:C:H4'	30:39:83:PHE:CD1	2.43	0.52
26:14:443:A:N7	30:39:45:ARG:HD2	2.24	0.52
46:H8:28:MET:HB2	46:H8:37:VAL:HG11	1.91	0.52
21:1F:2:GLY:O	21:1F:4:GLY:N	2.42	0.52
35:68:43:VAL:HG12	35:68:54:GLU:HA	1.90	0.52
26:14:2197:U:H1'	26:14:2198:A:C8	2.44	0.52
35:25:64:ARG:NH1	35:25:81:ASP:OD1	2.42	0.52
1:13:153:C:N4	1:13:168:G:H22	2.06	0.52
41:C8:106:PHE:HA	41:C8:109:LEU:HD12	1.90	0.52
31:41:47:LYS:NZ	31:41:80:PHE:HD1	2.07	0.52
7:6E:50:ILE:HB	7:6E:58:PRO:HB3	1.91	0.52
6:52:96:PRO:HB3	18:9A:30:ASP:CG	2.30	0.52
48:F5:52:ARG:HH11	48:F5:57:GLU:HG3	1.74	0.52
39:A8:34:HIS:HB2	39:A8:36:TYR:CE1	2.31	0.52
26:14:993:G:C5	26:14:994:C:H5	2.27	0.52
42:95:35:LEU:HB2	42:95:37:VAL:CG1	2.39	0.52
22:1K:46:7MG:C8	22:1K:46:7MG:H5''	2.43	0.52
29:21:119:ARG:HG3	29:21:119:ARG:NH1	2.24	0.52
1:13:1125:U:H5'	1:13:1126:U:C5	2.43	0.52
26:1H:963:U:OP1	58:1H:3799:HOH:O	2.19	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1G:1294:G:H2'	1:1G:1295:G:H8	1.74	0.52
9:82:103:THR:HG22	9:82:105:ASP:H	1.74	0.52
38:98:48:VAL:O	38:98:51:LEU:N	2.42	0.52
15:6I:56:LEU:O	15:6I:60:VAL:HG23	2.09	0.52
26:14:2346:A:H5''	26:14:2383:G:C1'	2.40	0.52
1:13:1051:C:H2'	1:13:1052:U:C6	2.43	0.52
36:78:38:GLN:HG2	36:78:45:LEU:HD12	1.92	0.52
1:1G:9:G:H1	1:1G:25:C:H42	1.57	0.52
1:1G:1308:U:H5''	13:4A:98:VAL:CG2	2.38	0.52
4:3E:161:ASN:O	4:3E:165:MET:HB2	2.09	0.52
28:11:25:THR:HB	28:11:82:ILE:H	1.74	0.52
41:C8:14:HIS:O	41:C8:18:LEU:HD12	2.08	0.52
27:16:44:G:H1'	27:16:47:C:N4	2.25	0.52
3:22:18:TRP:CD1	14:5A:54:PRO:HA	2.43	0.52
5:42:151:LEU:HD13	8:72:77:GLU:HG2	1.91	0.52
29:21:52:LEU:O	29:21:75:VAL:HG23	2.09	0.52
1:1G:1256:A:H62	1:1G:1277:C:H3'	1.72	0.52
26:14:567:A:OP2	36:35:29:LYS:NZ	2.42	0.52
1:1G:962:C:H42	1:1G:973:G:H1	1.57	0.52
36:78:49:ARG:HG3	55:Q8:58:ILE:HG13	1.92	0.52
26:1H:2392:A:H8	36:78:61:ARG:HG2	1.73	0.52
30:39:27:GLU:C	30:39:28:ILE:HG13	2.29	0.52
1:13:963:G:H5'	58:13:1866:HOH:O	2.08	0.52
1:13:1072:G:H2'	1:13:1073:U:O4'	2.10	0.52
48:J8:87:PRO:O	48:J8:91:LYS:HB2	2.10	0.52
26:14:2720:U:N3	26:14:2873:A:H2	2.06	0.52
1:13:631:G:C8	1:13:632:A:C2	2.98	0.52
24:3K:30:G:H1	24:3K:40:C:N4	2.07	0.52
2:12:196:LEU:HD12	2:12:197:VAL:HG23	1.91	0.52
9:82:119:ALA:O	9:82:120:ARG:HB2	2.10	0.52
38:98:79:LEU:HD23	38:98:83:ILE:HD12	1.91	0.52
26:1H:1324:G:C4	26:1H:1328:G:O6	2.62	0.52
26:14:1454:U:OP1	38:55:77:ARG:NH1	2.42	0.52
37:45:135:ASP:CG	46:D5:81:ARG:HH22	2.13	0.52
27:16:7:G:O5'	39:A8:29:PHE:CE2	2.63	0.52
7:62:26:PHE:CD2	7:62:30:ILE:HD11	2.45	0.52
52:N8:41:PRO:O	52:N8:44:THR:HG22	2.09	0.52
29:29:9:VAL:HG23	29:29:26:ILE:O	2.10	0.52
14:5I:48:ALA:HB2	14:5I:53:LEU:HD12	1.91	0.52
26:1H:1336:A:H2'	26:1H:1337:G:H8	1.73	0.52
1:1G:1340:A:O2'	23:2L:33:OMC:H5''	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1G:1122:U:N3	1:1G:1123:A:N7	2.58	0.52
37:88:11:LYS:HE2	37:88:86:GLY:HA2	1.90	0.52
26:1H:1268:A:H2'	26:1H:1269:A:O4'	2.09	0.52
26:14:2233:U:H2'	26:14:2234:G:C8	2.43	0.52
9:8E:77:ILE:O	9:8E:81:ILE:HG12	2.09	0.52
6:52:33:TYR:OH	6:52:78:GLU:HG3	2.09	0.52
26:1H:1110:G:O2'	26:1H:1111:A:O5'	2.25	0.52
1:1G:865:A:N3	1:1G:918:A:O2'	2.37	0.52
6:5E:39:LYS:HB2	6:5E:64:GLN:HB2	1.90	0.52
26:1H:1766:U:H2'	26:1H:1767:C:H6	1.74	0.52
47:18:64:ASP:HB2	47:18:85:ALA:HB1	1.90	0.52
33:69:124:GLY:H	33:69:142:VAL:HG12	1.74	0.52
1:1G:591:U:OP2	8:72:30:ARG:HD3	2.08	0.52
46:H8:126:VAL:HG12	46:H8:163:LEU:HD23	1.91	0.52
1:1G:998:G:H2'	1:1G:998(A):C:C6	2.44	0.52
4:32:134:ASP:O	4:32:136:PRO:HD3	2.09	0.52
26:14:2839:G:H21	38:55:92:GLY:HA2	1.73	0.52
1:1G:500:G:H2'	1:1G:501:C:C6	2.44	0.52
55:M5:9:GLY:O	55:M5:13:ARG:HD2	2.09	0.52
26:1H:2287:A:C2	26:1H:2289:G:C8	2.97	0.52
1:1G:862:C:H1'	1:1G:874:G:H5''	1.90	0.52
32:59:41:MET:N	32:59:41:MET:SD	2.82	0.52
26:14:2150:U:H2'	26:14:2151:G:C8	2.44	0.52
1:13:1060:C:O2'	10:1I:56:HIS:ND1	2.34	0.52
30:31:29:ASN:N	30:31:112:MET:HE1	2.25	0.52
28:19:17:THR:O	28:19:211:ARG:NH2	2.39	0.52
1:13:222:U:H2'	1:13:223:U:H6	1.74	0.52
26:1H:529:A:H8	26:1H:530:G:C6	2.26	0.52
6:52:23:LYS:NZ	6:52:42:GLU:OE2	2.30	0.52
26:1H:1126:A:H4'	26:1H:1127:A:O5'	2.09	0.52
2:1E:7:VAL:HG22	2:1E:8:LYS:H	1.75	0.52
1:13:1298:C:P	7:6E:114:ARG:HH22	2.32	0.52
26:14:2641:G:O3'	34:15:76:SER:OG	2.26	0.52
16:7I:13:HIS:C	16:7I:15:PRO:HD3	2.30	0.52
15:6I:7:GLU:OE1	15:6I:38:ARG:NH2	2.42	0.52
1:13:316:G:OP2	1:13:351:G:O2'	2.27	0.52
32:59:75:ALA:O	32:59:79:VAL:HG13	2.09	0.52
26:1H:2335:A:C8	26:1H:2337:G:C5	2.98	0.52
6:52:100:ASN:ND2	18:9A:26:LEU:O	2.41	0.52
5:4E:73:ASN:ND2	5:4E:73:ASN:O	2.40	0.52
26:14:548:A:C5	26:14:549:G:H1'	2.44	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:58:28:THR:N	34:58:106:MET:HE1	2.24	0.52
41:85:92:ARG:C	41:85:94:ASN:H	2.13	0.52
26:1H:2427:C:H5''	26:1H:2428:G:OP1	2.10	0.52
39:65:110:LEU:HB3	39:65:112:PHE:HE1	1.74	0.52
26:1H:1328:G:H2'	26:1H:1330:C:C5	2.45	0.52
1:1G:1096:C:O2'	1:1G:1170:A:O2'	2.21	0.52
26:1H:198:C:H5'	26:1H:2244:U:OP1	2.10	0.52
26:14:1058:U:H2'	26:14:1059:G:C8	2.45	0.52
20:BI:53:LEU:HD23	20:BI:100:ILE:HG22	1.92	0.52
20:BI:53:LEU:O	20:BI:57:ARG:HB2	2.10	0.52
53:O8:10:LEU:HD23	55:Q8:32:LEU:HD13	1.92	0.52
1:13:243:A:H4'	1:13:244:U:H5''	1.92	0.52
26:1H:1336:A:OP2	44:F8:64:LYS:NZ	2.34	0.52
1:1G:1411:C:H2'	1:1G:1412:C:C6	2.44	0.52
4:32:112:VAL:HG12	4:32:116:GLN:OE1	2.10	0.52
21:1B:2:GLY:O	21:1B:4:GLY:N	2.43	0.52
26:1H:773:U:C4'	28:11:47:GLY:HA3	2.39	0.52
2:12:145:LEU:O	2:12:149:LEU:HB2	2.10	0.52
2:1E:223:ILE:HG21	2:1E:230:VAL:HG22	1.90	0.52
26:14:99:U:H4'	26:14:101:G:H5'	1.90	0.52
1:13:984:C:H42	1:13:1221:G:H1	1.56	0.52
26:1H:2422:A:N7	55:Q8:30:ARG:NH2	2.55	0.52
37:45:103:MET:HB2	37:45:104:PHE:HD1	1.75	0.52
1:13:804:U:H5''	1:13:805:C:OP2	2.09	0.52
7:6E:5:ARG:NH2	7:6E:7:ALA:HA	2.25	0.52
26:14:2757:A:N1	32:59:67:LEU:HD22	2.24	0.52
45:G8:85:VAL:HG22	45:G8:98:VAL:HB	1.91	0.52
26:14:1111:A:H4'	32:59:3:ARG:HH11	1.73	0.52
26:14:1035:U:H2'	26:14:1036:G:C8	2.44	0.52
26:1H:654(B):C:H2'	26:1H:654(C):G:C8	2.44	0.52
55:M5:59:LYS:C	55:M5:60:LEU:HG	2.30	0.52
30:39:5:ALA:HB1	30:39:125:LEU:HD21	1.92	0.52
9:82:118:LYS:HB3	9:82:118:LYS:NZ	2.24	0.52
26:14:2776:A:H4'	26:14:2777:G:O5'	2.10	0.52
1:13:1142:G:H2'	1:13:1143:G:O4'	2.09	0.52
15:6I:74:ASP:OD1	15:6I:77:ARG:N	2.41	0.52
30:39:18:ARG:HG2	30:39:19:GLU:N	2.23	0.52
3:22:73:PRO:O	3:22:76:VAL:HG13	2.09	0.52
26:14:1796:U:H2'	26:14:1797:C:C6	2.45	0.52
37:45:31:ASP:H	37:45:107:ALA:HB2	1.73	0.52
26:1H:2179:C:H2'	26:1H:2180:U:H6	1.74	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:12:19:HIS:CG	2:12:20:GLU:H	2.28	0.52
41:85:66:ASN:OD1	41:85:76:TYR:N	2.42	0.52
37:88:85:LYS:HG3	37:88:86:GLY:N	2.24	0.52
26:1H:1111:A:H4'	32:51:3:ARG:HH11	1.74	0.52
26:14:2643:G:H2'	26:14:2644:G:O4'	2.09	0.52
34:15:33:LEU:HD12	34:15:38:HIS:CD2	2.45	0.52
6:52:49:ALA:HB2	18:9A:78:LEU:O	2.10	0.52
26:14:1309:G:N7	58:14:3924:HOH:O	2.34	0.52
11:2I:27:ASN:OD1	11:2I:28:THR:N	2.43	0.52
20:BA:67:ALA:O	20:BA:73:HIS:ND1	2.43	0.52
1:13:800:G:O6	58:13:1943:HOH:O	2.18	0.52
26:1H:1931:U:H5	26:1H:1969:A:N7	2.07	0.52
28:11:2:ALA:HA	28:11:20:ASP:HB2	1.91	0.52
26:14:2468:G:H3'	26:14:2476:A:N1	2.24	0.52
17:8I:81:ARG:NH2	17:8I:83:ASP:OD2	2.42	0.52
8:72:104:ARG:HB3	8:72:107:LEU:HB2	1.91	0.52
1:13:1502:A:H2	1:13:1505:G:N1	1.97	0.52
26:1H:1512:G:H2'	26:1H:1513:C:C6	2.45	0.52
20:BI:30:LYS:HZ3	20:BI:80:ARG:HH12	1.58	0.52
9:82:70:LYS:O	9:82:74:ILE:HG13	2.09	0.52
26:14:899:A:H2'	26:14:900:A:H8	1.73	0.52
26:1H:881:G:O6	26:1H:895:U:N3	2.42	0.52
46:D5:69:THR:HG22	46:D5:90:VAL:HG22	1.91	0.52
40:75:26:ASP:OD1	40:75:120:ARG:NH2	2.39	0.52
1:1G:872:A:O2'	1:1G:873:A:H5''	2.10	0.52
26:1H:66:C:H2'	26:1H:67:U:C6	2.45	0.52
26:1H:2801:A:H2'	26:1H:2802:G:H4'	1.90	0.52
26:14:944:G:O3'	58:14:3658:HOH:O	2.19	0.52
26:14:1386:C:OP2	26:14:1396:U:H5	1.92	0.52
26:1H:2123:G:H22	26:1H:2175:C:N4	2.08	0.52
32:59:30:LYS:HB3	32:59:79:VAL:HA	1.92	0.52
1:13:1441:G:H21	1:13:1460:A:H62	1.57	0.52
27:16:28:C:H2'	27:16:29:A:C8	2.45	0.52
26:1H:1443:G:C2	26:1H:1549:C:N3	2.77	0.52
26:14:2683:C:OP1	40:75:53:ARG:NH2	2.42	0.52
26:1H:533:G:H5'	41:C8:24:TYR:CE1	2.45	0.52
26:14:1131:G:O6	26:14:2040:C:H1'	2.10	0.52
15:6A:78:TYR:CZ	15:6A:82:ILE:HD12	2.45	0.52
20:BA:63:ILE:HG21	20:BA:81:LYS:HG3	1.92	0.52
34:15:61:ARG:NE	34:15:61:ARG:HA	2.25	0.52
9:8E:10:ARG:NE	9:8E:105:ASP:OD2	2.38	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:13:191(C):G:H2'	1:13:191(D):U:O4'	2.09	0.52
17:8A:67:LYS:O	17:8A:68:ARG:HB3	2.10	0.52
26:14:1810:A:H2'	26:14:1811:G:O4'	2.09	0.52
1:1G:16:A:H2'	1:1G:17:U:H6	1.75	0.52
26:14:2130:U:H2'	26:14:2158:A:C6	2.45	0.52
26:14:666:G:H5''	36:35:47:ASP:O	2.10	0.52
3:22:9:GLY:N	14:5A:49:HIS:O	2.43	0.52
26:14:2287:A:C2	26:14:2289:G:C8	2.98	0.52
30:31:67:GLN:O	30:31:67:GLN:HG3	2.08	0.52
26:14:1209:G:H21	26:14:1210:A:H62	1.56	0.52
50:H5:43:ILE:O	50:H5:47:VAL:HG23	2.10	0.52
29:29:11:MET:HE3	29:29:187:ALA:H	1.74	0.52
3:22:47:LEU:HD13	3:22:50:ALA:HB3	1.91	0.52
13:4I:27:LYS:HA	13:4I:31:LYS:NZ	2.25	0.52
1:1G:626:U:C2	1:1G:627:G:C8	2.98	0.52
53:O8:25:LYS:HB2	55:Q8:32:LEU:HD12	1.91	0.52
37:45:26:TYR:O	37:45:26:TYR:CD1	2.63	0.52
26:1H:2109:U:N3	26:1H:2110:G:O6	2.41	0.52
1:1G:866:C:O2'	1:1G:919:A:OP1	2.23	0.52
1:13:1352:C:OP1	21:1F:3:LYS:NZ	2.29	0.52
8:7E:64:LYS:HG2	8:7E:79:VAL:HG21	1.91	0.52
26:1H:2846:G:N7	58:1H:4384:HOH:O	2.34	0.52
26:1H:455:C:N3	26:1H:473:G:H5'	2.25	0.52
26:14:1028:A:N3	26:14:2486:G:O2'	2.34	0.52
26:1H:2785:C:H2'	26:1H:2786:U:O4'	2.09	0.52
49:K8:24:LEU:HD13	49:K8:60:LEU:HD11	1.91	0.52
6:5E:41:GLU:HB2	6:5E:62:TRP:CE3	2.45	0.52
24:1L:30:G:H1	24:1L:40:C:H42	1.57	0.52
1:1G:807:A:H2'	1:1G:808:C:C6	2.44	0.52
26:14:71:A:H5'	26:14:71:A:H8	1.70	0.52
1:13:520:A:N1	1:13:536:C:H1'	2.25	0.52
1:1G:1435:G:H2'	1:1G:1436:U:H6	1.74	0.52
30:39:117:ARG:NH1	30:39:120:GLU:OE1	2.43	0.52
26:14:2068:U:N3	26:14:2430:A:H2	2.03	0.52
48:J8:91:LYS:HA	48:J8:91:LYS:HZ3	1.74	0.52
26:1H:1164:G:H2'	26:1H:1165:U:C6	2.45	0.52
2:1E:210:SER:O	2:1E:214:ILE:HG12	2.09	0.52
26:14:1533:C:N3	26:14:1534:G:O2'	2.39	0.52
31:41:173:LEU:HB3	31:41:178:PHE:CD2	2.43	0.52
26:1H:2232:U:OP1	48:J8:40:ARG:NH1	2.42	0.52
1:1G:538:G:OP2	12:3A:115:LYS:HB2	2.10	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:13:1321:C:H3'	1:13:1322:C:H5''	1.92	0.52
1:13:953:G:H2'	1:13:954:G:O4'	2.10	0.52
26:1H:270(L):U:O2	33:61:50:ARG:HG2	2.09	0.52
26:1H:547:A:C6	26:1H:548:A:C6	2.97	0.52
1:1G:1328:C:O2'	13:4A:29:ARG:NE	2.41	0.52
26:14:1057:A:H2'	26:14:1058:U:O4'	2.09	0.52
26:1H:2690:C:H5''	26:1H:2872:G:N2	2.24	0.52
41:C8:50:ARG:NH1	42:D8:72:VAL:HG12	2.25	0.52
26:1H:1486:A:H2'	26:1H:1487:G:C8	2.45	0.52
32:59:77:LYS:HD2	32:59:80:SER:HB2	1.90	0.52
26:1H:274:G:H2'	26:1H:275:G:H4'	1.92	0.52
26:14:118:A:N3	26:14:178:G:H1'	2.25	0.52
40:75:61:PHE:CE1	40:75:76:PHE:HB2	2.44	0.52
48:F5:29:GLY:O	48:F5:30:VAL:HG22	2.10	0.52
8:72:16:ALA:HA	8:72:19:VAL:HG22	1.91	0.52
31:41:166:ASP:N	31:41:166:ASP:OD1	2.42	0.52
32:51:92:ILE:HD11	32:51:160:LYS:HZ2	1.75	0.52
36:78:130:PHE:HE1	36:78:146:VAL:HG23	1.75	0.52
26:1H:1262:A:N3	52:N8:10:LYS:HE3	2.25	0.52
26:14:1451:C:H3'	26:14:1453:A:H5'	1.90	0.52
29:21:135:HIS:CE1	58:21:402:HOH:O	2.58	0.52
24:3L:36:A:N7	25:4L:14:A:N6	2.58	0.52
26:14:1142(A):A:H4'	34:15:25:ARG:HH22	1.75	0.52
55:Q8:54:GLU:O	55:Q8:56:GLU:N	2.42	0.52
1:13:558:G:H2'	1:13:559:A:H2	1.74	0.52
26:1H:2056:G:C2	26:1H:2057:A:C8	2.98	0.52
1:1G:1046:A:H61	1:1G:1213:A:H61	1.58	0.52
1:1G:164:U:H2'	1:1G:165:C:C6	2.44	0.52
2:12:162:ILE:HD11	2:12:184:VAL:HG13	1.91	0.52
31:41:67:LYS:HZ3	51:M8:6:HIS:CE1	2.28	0.52
13:4A:37:THR:O	13:4A:55:ARG:NE	2.43	0.52
1:1G:1259:C:O2	1:1G:1283:G:O2'	2.27	0.52
27:16:7:G:H5''	27:16:7:G:H8	1.75	0.52
39:A8:29:PHE:HD1	39:A8:30:ARG:N	2.08	0.52
1:13:232:G:H2'	1:13:233:C:H6	1.76	0.52
26:14:322:A:OP2	30:39:169:ASN:HB2	2.10	0.52
28:19:16:MET:HE1	28:19:208:LYS:HG2	1.91	0.52
43:E8:78:GLU:OE1	43:E8:99:ARG:HD3	2.10	0.52
4:3E:108:LEU:HB3	4:3E:110:PHE:CE1	2.45	0.52
31:41:7:LEU:HD11	31:41:176:LEU:HD22	1.91	0.52
26:14:1847:A:O2'	26:14:1848:A:H5'	2.10	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
12:3A:86:ARG:HB2	12:3A:101:VAL:HG23	1.91	0.52
26:1H:1190:G:C5	58:1H:3819:HOH:O	2.63	0.52
1:1G:278:G:OP2	17:8A:41:LYS:NZ	2.34	0.52
38:55:12:ARG:HD3	38:55:16:HIS:CD2	2.45	0.52
49:K8:33:MET:O	49:K8:37:PHE:HD1	1.93	0.52
26:14:1903:G:OP1	28:19:241:PRO:HB2	2.09	0.52
49:K8:14:ARG:NH1	49:K8:66:GLU:OE2	2.43	0.52
1:13:1348:U:H4'	9:8E:120:ARG:HD2	1.91	0.51
1:1G:975:A:H4'	1:1G:976:G:C5'	2.41	0.51
1:13:767:A:H2'	1:13:768:A:O4'	2.09	0.51
24:3K:13:C:H2'	24:3K:14:A:C8	2.33	0.51
1:1G:20:U:H2'	1:1G:21:G:O4'	2.10	0.51
5:42:5:ASP:HA	5:42:63:ARG:NH1	2.25	0.51
1:13:323:U:H2'	1:13:324:G:O4'	2.10	0.51
29:21:39:PRO:HD3	29:21:45:THR:HG22	1.92	0.51
6:52:97:PHE:O	18:9A:31:LEU:HD23	2.10	0.51
1:13:148:G:H2'	1:13:149:A:H8	1.75	0.51
35:68:120:GLU:HG2	35:68:122:LEU:HG	1.92	0.51
26:14:2557:G:H2'	26:14:2558:C:H6	1.73	0.51
39:65:62:LYS:HA	39:65:65:VAL:HG12	1.92	0.51
45:G8:28:LYS:HD2	45:G8:40:GLU:OE2	2.09	0.51
29:29:120:TRP:CE3	29:29:155:LYS:HD3	2.45	0.51
26:14:2808:U:H5''	26:14:2891:G:O6	2.10	0.51
13:4I:52:GLU:O	13:4I:56:LEU:HB2	2.09	0.51
9:82:48:GLU:OE1	9:82:51:ARG:NH1	2.43	0.51
51:I5:48:ARG:NH1	51:I5:51:ASP:HB3	2.25	0.51
7:6E:69:VAL:HG22	7:6E:135:VAL:HG22	1.92	0.51
26:1H:747:U:O2	26:1H:2014:A:H1'	2.10	0.51
26:14:1599:C:H2'	26:14:1600:C:H6	1.75	0.51
16:7I:58:TYR:O	16:7I:62:VAL:HG22	2.09	0.51
26:14:1322:A:N1	26:14:1333:C:O2'	2.35	0.51
38:55:45:ARG:HA	38:55:95:THR:HG21	1.93	0.51
24:3K:37:A:H2'	24:3K:38:A:O4'	2.11	0.51
41:C8:52:ARG:HA	41:C8:55:ARG:HG3	1.93	0.51
26:14:1416:G:H1	26:14:1582:C:N4	1.98	0.51
1:1G:1256:A:N6	1:1G:1277:C:H3'	2.25	0.51
1:1G:590:C:H2'	1:1G:591:U:C6	2.45	0.51
26:14:273(D):C:H42	26:14:363(B):G:H1	1.56	0.51
32:51:83:TYR:CB	32:51:134:SER:HA	2.39	0.51
2:12:174:VAL:HG11	2:12:196:LEU:HD13	1.90	0.51
40:75:4:GLY:N	40:75:7:ILE:HG22	2.25	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:1H:2157:G:O2'	26:1H:2158:A:O5'	2.25	0.51
1:13:1131:G:H2'	1:13:1132:C:C6	2.44	0.51
1:1G:1028:C:H2'	1:1G:1028(A):C:O4'	2.10	0.51
1:1G:1181:G:N7	1:1G:1182:G:N2	2.54	0.51
26:14:172:C:H2'	26:14:173:G:C8	2.44	0.51
41:C8:17:ILE:HD12	41:C8:32:PHE:HE1	1.75	0.51
27:16:44:G:C2	27:16:48:A:C2	2.97	0.51
26:14:2850:A:C2	26:14:2851:A:C4	2.98	0.51
26:14:2037:G:H2'	26:14:2038:G:C8	2.45	0.51
37:45:91:GLU:HG3	37:45:92:GLY:N	2.25	0.51
13:4A:102:ARG:HH11	13:4A:105:THR:HG23	1.75	0.51
33:69:7:GLU:HG3	33:69:8:PRO:HD2	1.92	0.51
1:1G:1387:G:H2'	1:1G:1388:C:C6	2.45	0.51
27:16:5:C:O2'	27:16:27:C:O2	2.27	0.51
50:H5:10:LYS:NZ	50:H5:15:TYR:OH	2.43	0.51
24:1L:38:A:H2'	24:1L:39:U:C6	2.45	0.51
23:2K:34:U:N3	23:2K:37:U:OP2	2.40	0.51
26:14:1259:G:H2'	26:14:1260:G:C8	2.45	0.51
26:1H:1827:C:H3'	58:1H:3705:HOH:O	2.10	0.51
26:1H:141(A):C:H2'	26:1H:142:G:O4'	2.09	0.51
27:1J:117:G:H8	27:1J:117:G:O5'	1.93	0.51
26:14:2749:A:H5''	32:59:6:ARG:NH2	2.26	0.51
26:14:1686:C:H2'	26:14:1687:G:O4'	2.10	0.51
40:B8:3:ARG:O	40:B8:7:ILE:N	2.43	0.51
55:M5:22:VAL:HG12	55:M5:50:LEU:HD23	1.91	0.51
1:1G:458:C:H2'	1:1G:464:G:C8	2.45	0.51
26:1H:412:A:H5''	26:1H:413:C:OP2	2.11	0.51
26:1H:1800:C:OP2	28:11:183:ARG:NH2	2.43	0.51
26:1H:579:G:H2'	26:1H:580:C:C6	2.46	0.51
26:1H:1103:A:H3'	26:1H:1104:C:C6	2.45	0.51
26:14:2122:U:H2'	26:14:2123:G:O4'	2.09	0.51
18:9A:53:ARG:HA	18:9A:56:THR:OG1	2.10	0.51
33:61:68:LEU:HA	33:61:71:ILE:CG2	2.40	0.51
20:BI:57:ARG:NH1	20:BI:102:GLY:HA2	2.23	0.51
2:12:19:HIS:CE1	2:12:206:ASP:H	2.29	0.51
39:65:26:LEU:HD22	39:65:87:PHE:CE1	2.45	0.51
1:1G:313:A:H2'	1:1G:314:C:H6	1.75	0.51
39:65:66:ALA:HA	39:65:69:VAL:HG12	1.93	0.51
39:A8:103:GLU:O	39:A8:106:ARG:HD3	2.11	0.51
26:14:2468:G:H3'	26:14:2476:A:C2	2.45	0.51
37:45:37:LEU:HB2	37:45:128:LYS:O	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:1H:2564:A:C2	26:1H:2647:U:H4'	2.45	0.51
1:1G:921:U:O2	5:42:19:MET:HB3	2.11	0.51
7:62:47:CYS:O	7:62:50:ILE:HB	2.09	0.51
1:1G:56:U:H2'	1:1G:57:G:C8	2.45	0.51
1:1G:1074:G:O2'	1:1G:1101:A:N1	2.34	0.51
26:1H:602:G:O2'	26:1H:655:A:N6	2.43	0.51
26:1H:6:A:N3	34:58:131:GLN:HG3	2.24	0.51
3:2E:18:TRP:H	3:2E:18:TRP:HE3	1.58	0.51
26:14:2019:A:H4'	41:85:34:LYS:HD2	1.92	0.51
30:39:158:THR:HB	30:39:195:ASP:HB2	1.92	0.51
4:32:30:LYS:CB	4:32:35:ARG:HD2	2.39	0.51
9:8E:24:GLY:HA2	9:8E:59:PHE:O	2.10	0.51
34:15:4:TYR:CD2	41:85:100:VAL:HG11	2.46	0.51
1:13:701:C:O2	1:13:703:G:N1	2.44	0.51
26:1H:880:G:H2'	26:1H:881:G:C8	2.46	0.51
26:1H:1312:U:H4'	26:1H:1313:U:O5'	2.10	0.51
44:F8:27:THR:HB	44:F8:80:ILE:HG22	1.91	0.51
26:1H:443:A:N7	30:31:45:ARG:HG2	2.26	0.51
29:29:25:VAL:HG12	29:29:26:ILE:H	1.74	0.51
26:14:1027:A:H5'	27:1J:88:C:H41	1.75	0.51
26:1H:2690:C:H5''	26:1H:2872:G:H21	1.74	0.51
1:13:1318:A:H1'	19:AI:37:ARG:HH21	1.75	0.51
1:1G:78:G:N2	1:1G:91:C:O2	2.42	0.51
26:1H:1026:U:H1'	26:1H:1027:A:P	2.51	0.51
2:1E:7:VAL:HB	2:1E:217:ARG:HD2	1.91	0.51
1:13:1313:U:O4	19:AI:4:SER:HB2	2.10	0.51
7:62:50:ILE:HG21	7:62:58:PRO:HA	1.91	0.51
26:1H:7:G:N2	26:1H:8:A:H1'	2.26	0.51
1:1G:179:A:H2'	1:1G:180:U:H6	1.74	0.51
1:13:1478:C:H2'	1:13:1479:C:C6	2.45	0.51
1:13:1442:G:C6	1:13:1446:A:C6	2.99	0.51
1:13:4:U:C4	8:7E:102:ARG:HG3	2.45	0.51
5:42:24:ARG:HB2	5:42:26:PHE:CE2	2.45	0.51
26:1H:1444:G:C2	26:1H:1548:C:N3	2.79	0.51
1:13:1117:G:OP1	9:8E:9:ARG:NH2	2.39	0.51
26:14:1014:U:N3	26:14:1015:G:N7	2.58	0.51
44:B5:25:LYS:HA	44:B5:81:VAL:O	2.10	0.51
17:8A:45:HIS:O	17:8A:73:VAL:HG12	2.10	0.51
6:5E:98:LEU:HA	18:9I:29:PHE:O	2.10	0.51
11:2I:34:ASP:HB3	11:2I:40:ILE:HD11	1.91	0.51
1:1G:243:A:H4'	1:1G:244:U:O5'	2.09	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:42:145:LYS:HA	5:42:148:VAL:HB	1.93	0.51
1:1G:1148:U:H2'	1:1G:1149:C:O4'	2.11	0.51
1:13:1374:A:O2'	7:6E:28:ASN:HB3	2.11	0.51
4:32:53:ASP:OD2	5:42:107:ARG:NH2	2.43	0.51
24:3K:15:G:H1	24:3K:48:C:N4	2.04	0.51
31:49:95:ARG:HG2	31:49:96:ARG:HG3	1.91	0.51
39:65:78:LEU:HD11	39:65:107:GLU:HB3	1.93	0.51
33:61:40:THR:H	33:61:43:ASN:HB2	1.75	0.51
26:1H:2250:G:C8	26:1H:2496:C:H5''	2.45	0.51
26:1H:882:G:H22	26:1H:894:C:H42	1.58	0.51
1:1G:502:G:C2	1:1G:503:C:C2	2.98	0.51
26:1H:1826:G:H4'	28:11:242:ARG:CZ	2.41	0.51
36:78:15:ARG:HA	36:78:16:ARG:HB2	1.93	0.51
26:1H:2636:U:OP1	29:21:80:GLU:HG3	2.09	0.51
1:1G:1315:U:H2'	1:1G:1316:G:O4'	2.11	0.51
31:41:35:GLU:HG3	31:41:36:LYS:HB3	1.93	0.51
26:1H:779:U:O4	58:1H:4240:HOH:O	2.19	0.51
20:BI:53:LEU:HA	20:BI:56:MET:HB3	1.93	0.51
26:1H:1027:A:C2	26:1H:2488:A:H5'	2.45	0.51
27:1J:94:C:H2'	27:1J:95:U:H6	1.76	0.51
27:1J:99:A:H3'	58:1J:313:HOH:O	2.09	0.51
26:1H:2001:A:H2'	26:1H:2002:G:C8	2.46	0.51
1:1G:993:G:O2'	1:1G:994:A:N7	2.40	0.51
26:14:1936:A:C8	26:14:1940:U:O2	2.63	0.51
1:1G:107:G:C2	1:1G:108:G:H1'	2.45	0.51
50:L8:12:PRO:O	50:L8:20:LYS:NZ	2.43	0.51
26:1H:1248:G:N2	30:31:88:VAL:HG21	2.26	0.51
1:1G:162:A:O5'	1:1G:162:A:H8	1.93	0.51
1:13:1278:U:H5''	1:13:1279:A:O4'	2.11	0.51
23:2L:44:A:H2'	23:2L:45:A:C8	2.45	0.51
36:35:15:ARG:NH2	36:35:17:LYS:HE3	2.24	0.51
26:14:993:G:C5	26:14:994:C:C5	2.99	0.51
22:1K:76:A:H8	26:1H:2583:G:H21	1.59	0.51
29:29:169:ASN:OD1	29:29:203:LYS:HB3	2.10	0.51
22:1K:22:G:H2'	22:1K:23:A:C8	2.45	0.51
1:1G:1048:G:O4'	1:1G:1215:G:H4'	2.10	0.51
1:13:660:G:H2'	1:13:661:G:H8	1.74	0.51
1:13:130:A:OP2	17:8I:63:ARG:NH2	2.42	0.51
26:14:1993:U:H4'	29:29:128:SER:HB3	1.91	0.51
10:1A:8:LEU:HB3	10:1A:16:LEU:HD22	1.92	0.51
12:3A:36:VAL:O	12:3A:59:ARG:N	2.44	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
37:45:74:TYR:O	37:45:89:ASN:HB2	2.11	0.51
5:4E:77:PRO:HG2	5:4E:142:LEU:HD22	1.93	0.51
1:1G:1130:A:H62	1:1G:1144:G:H21	1.59	0.51
23:2K:65:G:H2'	23:2K:66:C:O4'	2.11	0.51
17:8I:11:VAL:HG12	17:8I:85:VAL:HG22	1.93	0.51
34:15:128:HIS:CE1	34:15:130:HIS:HB3	2.46	0.51
26:1H:303:U:H2'	26:1H:304:G:H8	1.76	0.51
26:1H:950:G:H2'	26:1H:951:C:C6	2.46	0.51
5:42:105:VAL:HG21	5:42:128:PRO:HB3	1.92	0.51
5:4E:122:GLU:HG2	5:4E:131:ILE:HD12	1.92	0.51
11:2I:123:LYS:O	11:2I:126:ARG:HG3	2.11	0.51
44:B5:36:LYS:HG2	44:B5:54:VAL:HB	1.93	0.51
1:1G:1502:A:H2	1:1G:1505:G:N1	2.00	0.51
26:14:1110:G:O2'	32:59:3:ARG:NH1	2.44	0.51
1:13:60:A:N6	1:13:110:C:N3	2.57	0.51
1:1G:1257:U:H5'	1:1G:1258:G:C8	2.46	0.51
29:29:57:LYS:H	29:29:57:LYS:NZ	2.09	0.51
38:98:63:ARG:HB2	38:98:80:PHE:HE2	1.75	0.51
29:29:51:PHE:O	29:29:52:LEU:HB2	2.09	0.51
26:1H:1243:G:H4'	36:78:7:ARG:HH21	1.76	0.51
33:69:75:LEU:HD13	33:69:76:THR:N	2.25	0.51
26:14:1053:C:H2'	26:14:1054:A:O4'	2.11	0.51
41:85:85:LYS:HB3	41:85:116:ALA:HB1	1.93	0.51
26:1H:723:G:H2'	26:1H:724:U:O4'	2.10	0.51
45:G8:5:MET:HE1	45:G8:32:PRO:HA	1.92	0.51
1:13:1410:G:H2'	1:13:1411:C:H6	1.76	0.51
6:52:12:PRO:HB3	6:52:58:GLY:HA2	1.93	0.51
5:42:33:VAL:HG21	5:42:109:ILE:HG12	1.93	0.51
26:14:1849:G:H2'	26:14:1850:G:H8	1.74	0.51
9:82:89:ASN:O	9:82:92:TYR:HB2	2.10	0.51
1:13:964:A:N3	1:13:969:A:O2'	2.32	0.51
2:1E:94:ASN:OD1	2:1E:95:GLN:N	2.40	0.51
18:9I:25:THR:HB	18:9I:42:ARG:HH12	1.75	0.51
1:13:240:C:H2'	1:13:241:C:C6	2.46	0.51
54:L5:12:ARG:HH21	54:L5:44:PRO:HB3	1.75	0.51
1:13:717:C:H6	1:13:717:C:H5''	1.76	0.51
42:D8:45:THR:OG1	42:D8:45:THR:O	2.24	0.51
43:E8:57:ASN:HA	43:E8:61:ASN:HD22	1.74	0.51
26:1H:1520:U:H2'	26:1H:1521:G:O4'	2.11	0.51
1:13:407:G:O4'	4:3E:119:GLN:NE2	2.43	0.51
38:98:3:HIS:O	38:98:5:LYS:N	2.44	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:802:A:H4'	58:14:3855:HOH:O	2.10	0.51
2:12:5:ILE:HA	2:12:221:LEU:HD21	1.92	0.51
55:Q8:49:VAL:CG1	55:Q8:52:LYS:HG3	2.39	0.51
45:C5:14:LEU:HD12	45:C5:15:VAL:N	2.25	0.51
1:1G:617:G:H1	1:1G:623:C:N4	2.04	0.51
26:14:1856:G:N2	26:14:1886:C:N3	2.48	0.51
16:7I:79:VAL:HG12	16:7I:80:PHE:CD1	2.45	0.51
34:15:42:TRP:O	41:85:64:ARG:HD2	2.10	0.51
2:12:178:ARG:HH11	2:12:178:ARG:HB2	1.75	0.51
26:14:481:G:OP2	45:C5:47:LYS:HB2	2.10	0.51
26:14:2310:A:H5'	26:14:2311:A:OP2	2.11	0.51
46:D5:126:VAL:HA	46:D5:163:LEU:HA	1.93	0.51
1:13:547:A:OP2	4:3E:2:GLY:HA2	2.11	0.51
26:14:747:U:OP2	52:J5:3:LYS:HD3	2.11	0.51
49:K8:42:GLY:C	49:K8:44:LEU:H	2.14	0.51
1:1G:1227:A:N3	19:AA:84:GLY:HA2	2.25	0.51
26:14:2105:C:H2'	26:14:2106:G:O4'	2.10	0.51
1:13:1417:G:N2	1:13:1482:G:H2'	2.25	0.51
20:BA:51:GLU:HA	20:BA:54:LYS:HB3	1.93	0.51
26:1H:934:G:H2'	26:1H:935:C:H6	1.76	0.51
26:1H:1204:A:H2	26:1H:1241:A:N1	2.08	0.51
46:H8:52:SER:O	46:H8:52:SER:OG	2.21	0.51
29:29:37:ARG:HD3	29:29:44:TYR:CE1	2.46	0.51
45:G8:93:GLY:O	45:G8:94:LYS:HB2	2.11	0.51
29:21:147:PRO:HB2	29:21:149:ARG:HG3	1.93	0.51
55:Q8:49:VAL:HG22	55:Q8:50:LEU:N	2.21	0.51
1:1G:1062:U:H2'	1:1G:1063:C:C6	2.45	0.51
37:88:135:ASP:HB3	37:88:137:TYR:N	2.26	0.51
30:39:148:LEU:HD21	30:39:191:ARG:HH11	1.75	0.51
28:19:43:ARG:HD2	28:19:49:ILE:HB	1.91	0.51
23:2L:24:C:C2	23:2L:25:U:C5	2.99	0.51
1:1G:1096:C:HO2'	1:1G:1170:A:HO2'	1.48	0.51
33:61:62:LYS:HG3	33:61:133:HIS:NE2	2.26	0.51
33:69:77:LEU:HD12	33:69:78:THR:H	1.76	0.51
1:1G:1411:C:H2'	1:1G:1412:C:H6	1.76	0.51
37:88:65:PHE:O	37:88:66:ILE:HB	2.11	0.51
1:13:1410:G:H2'	1:13:1411:C:C6	2.45	0.51
43:A5:14:PRO:HG2	43:A5:78:GLU:HG3	1.92	0.51
1:1G:35:G:C2	1:1G:550:G:N3	2.79	0.51
26:14:380:U:H5'	48:F5:18:ILE:HD12	1.93	0.51
26:14:1448:G:H1'	26:14:1528:A:H62	1.76	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
12:3A:53:ARG:HH12	12:3A:92:ASP:HB3	1.75	0.51
26:14:1842:G:O2'	28:19:253:GLN:OE1	2.28	0.51
1:1G:15:G:N7	1:1G:1396:A:C2	2.79	0.51
1:1G:858:G:H8	1:1G:858:G:OP2	1.94	0.51
26:1H:754:C:H2'	26:1H:755:C:C6	2.46	0.51
26:14:907:U:O5'	37:45:24:GLY:HA2	2.11	0.51
26:1H:821:A:H5''	26:1H:822:U:H6	1.76	0.51
26:14:1336:A:H2'	26:14:1337:G:C8	2.46	0.51
33:61:127:VAL:HA	33:61:138:ILE:O	2.11	0.51
37:45:112:GLU:HA	37:45:115:MET:HB2	1.93	0.51
1:1G:1119:C:OP2	9:82:9:ARG:NH2	2.41	0.51
26:1H:2394:C:H2'	26:1H:2395:C:H6	1.76	0.51
26:1H:1187:G:H5''	42:D8:81:TYR:CE1	2.46	0.51
1:1G:1392:G:H21	1:1G:1502:A:H8	1.57	0.51
47:E5:18:ALA:HB3	47:E5:20:ARG:HE	1.76	0.51
26:14:1021:A:H8	26:14:1021:A:H3'	1.75	0.51
26:1H:2160:G:C2	26:1H:2161:C:H1'	2.46	0.51
29:29:103:ASP:OD1	29:29:201:THR:HG23	2.11	0.51
26:1H:780:G:H21	26:1H:783:A:N6	2.01	0.51
1:13:1178:G:N2	1:13:1181:G:H8	2.09	0.51
27:1J:40:U:N3	27:1J:43:C:H5''	2.26	0.51
26:1H:2199:A:H5''	26:1H:2205:C:C5	2.41	0.51
1:1G:1060:C:HO2'	10:1A:56:HIS:HD1	1.57	0.51
1:1G:838:G:N2	1:1G:849:C:N3	2.58	0.51
26:14:2295:C:P	39:65:10:ARG:HD3	2.50	0.51
36:78:1:MET:HE2	36:78:6:LEU:HD13	1.92	0.51
26:1H:583:G:N7	58:1H:3992:HOH:O	2.35	0.51
2:1E:172:ILE:O	2:1E:176:GLU:HG3	2.11	0.51
1:1G:668:G:O4'	15:6A:49:ASP:HB2	2.11	0.51
23:2K:22:A:H8	23:2K:22:A:H5''	1.76	0.51
32:51:86:GLU:HG3	32:51:165:ALA:HB3	1.92	0.51
26:14:2257:U:O2'	26:14:2258:C:H5'	2.11	0.51
26:14:2104:G:H2'	26:14:2105:C:C6	2.46	0.51
1:1G:452:A:H62	1:1G:480:U:H3	1.59	0.51
1:1G:179:A:H2'	1:1G:180:U:C6	2.45	0.51
51:I5:49:PHE:HD2	51:I5:50:VAL:HG22	1.75	0.51
1:13:105:G:H2'	1:13:106:C:C6	2.46	0.51
20:BA:26:ASN:HA	20:BA:29:LYS:HG2	1.92	0.51
19:AA:29:ARG:NH1	19:AA:48:THR:HG23	2.26	0.51
21:1F:9:ARG:O	21:1F:13:ILE:HG13	2.11	0.51
45:C5:48:ALA:HB3	45:C5:59:GLY:C	2.31	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
38:55:56:LYS:NZ	38:55:90:ARG:O	2.42	0.51
31:49:106:LEU:HA	31:49:110:ALA:HB3	1.91	0.51
31:49:107:LEU:HD21	31:49:178:PHE:CE2	2.46	0.51
26:14:64:A:C4	44:B5:66:LEU:HD12	2.45	0.51
45:C5:62:GLU:CD	45:C5:63:LYS:H	2.15	0.51
11:2I:20:TYR:HB2	11:2I:31:THR:HG23	1.91	0.51
39:A8:78:LEU:HD12	39:A8:108:GLY:HA2	1.91	0.51
11:2I:57:THR:HG22	11:2I:59:TYR:H	1.75	0.51
26:14:2431:U:O2'	26:14:2433:A:N7	2.38	0.51
47:I8:51:VAL:N	47:I8:62:LEU:HD12	2.25	0.51
26:14:1654:A:H1'	26:14:2823:A:H5'	1.93	0.51
26:14:648:G:O2'	26:14:2351:G:OP1	2.20	0.51
26:1H:1427:A:H4'	26:1H:1428:C:O5'	2.10	0.51
29:21:4:ILE:HD13	29:21:28:ALA:HB1	1.92	0.51
30:31:101:LEU:HD13	30:31:102:PRO:HD2	1.93	0.50
1:13:1349:A:OP2	9:8E:118:LYS:NZ	2.36	0.50
26:1H:2702:U:C6	26:1H:2702:U:OP1	2.64	0.50
26:1H:453:C:OP1	58:1H:3841:HOH:O	2.19	0.50
1:13:1199:U:H4'	10:1I:54:PHE:CE2	2.47	0.50
26:1H:1242:A:N1	36:78:4:SER:OG	2.43	0.50
29:29:81:ILE:HG22	29:29:82:ARG:N	2.22	0.50
1:13:827:U:C5	1:13:872:A:N1	2.74	0.50
55:M5:14:VAL:CG1	55:M5:22:VAL:HG13	2.42	0.50
4:32:202:LEU:O	4:32:206:PHE:N	2.45	0.50
26:1H:2695:C:H2'	26:1H:2696:U:C6	2.43	0.50
24:3L:76:A:O2'	26:14:2394:C:N3	2.39	0.50
29:29:57:LYS:HZ2	29:29:57:LYS:H	1.58	0.50
36:78:6:LEU:O	36:78:7:ARG:HG2	2.12	0.50
4:32:108:LEU:HD21	4:32:183:GLY:HA3	1.94	0.50
26:1H:2531:A:H5'	32:51:157:TYR:CE1	2.45	0.50
26:1H:2734:A:H3'	26:1H:2735:G:H8	1.75	0.50
26:1H:536:A:H2'	26:1H:537:C:C6	2.46	0.50
3:2E:149:ALA:HA	3:2E:201:TYR:O	2.11	0.50
26:1H:1443:G:N2	26:1H:1549:C:N3	2.58	0.50
45:C5:43:ASN:HB3	45:C5:64:GLU:HA	1.93	0.50
39:65:64:GLU:O	39:65:68:GLN:HG3	2.11	0.50
26:14:2697:G:H2'	26:14:2698:U:O4'	2.10	0.50
26:1H:44:A:O2'	26:1H:45:G:H5'	2.11	0.50
28:19:166:GLN:HB3	28:19:174:ILE:HG22	1.92	0.50
26:14:77:C:OP1	49:G5:59:ARG:HD3	2.11	0.50
26:1H:1290:C:H2'	26:1H:1291:C:C6	2.46	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:1H:2408:U:H2'	26:1H:2409:G:C8	2.47	0.50
26:14:400:G:O6	58:14:3683:HOH:O	2.20	0.50
8:7E:16:ALA:HB1	8:7E:24:THR:HG21	1.93	0.50
4:3E:150:GLU:N	4:3E:150:GLU:OE1	2.45	0.50
30:39:113:ALA:HB1	30:39:186:ILE:HG21	1.93	0.50
11:2I:17:GLY:HA3	11:2I:77:MET:SD	2.51	0.50
27:1J:13:A:H5''	27:1J:15:A:N6	2.25	0.50
1:1G:976:G:H5'	1:1G:1358:U:O2'	2.11	0.50
24:3K:18:G:HO2'	24:3K:19:G:P	2.34	0.50
26:1H:2199:A:H3'	26:1H:2205:C:H6	1.76	0.50
26:1H:1668:A:H61	26:1H:1676:A:H61	1.58	0.50
28:19:43:ARG:HA	28:19:49:ILE:HA	1.93	0.50
3:2E:52:LEU:HA	3:2E:70:VAL:HG12	1.93	0.50
31:49:53:LEU:HD21	31:49:87:PRO:HB2	1.94	0.50
26:14:1277:G:O2'	38:55:24:GLN:HG2	2.11	0.50
2:12:166:ASP:OD1	2:12:168:THR:OG1	2.28	0.50
1:1G:222:U:H2'	1:1G:223:U:H6	1.74	0.50
26:1H:1102:C:H2'	26:1H:1103:A:C8	2.46	0.50
10:1A:99:LYS:CD	10:1A:100:THR:H	2.23	0.50
1:13:674:G:H2'	1:13:675:A:C8	2.45	0.50
26:14:864:G:C6	26:14:865:C:N4	2.79	0.50
1:13:153:C:H42	1:13:168:G:H22	1.58	0.50
24:3L:17:C:N4	26:14:2112:G:OP1	2.44	0.50
31:41:110:ALA:HA	31:41:140:ILE:O	2.11	0.50
18:9A:74:ARG:HB3	18:9A:81:PHE:CE1	2.46	0.50
1:1G:1469:G:H2'	1:1G:1470:G:H8	1.76	0.50
4:3E:43:HIS:O	4:3E:46:LYS:HG2	2.12	0.50
39:A8:66:ALA:HA	39:A8:69:VAL:HG12	1.93	0.50
29:21:15:PHE:HB3	40:B8:81:PRO:HG3	1.92	0.50
1:1G:79:G:H2'	1:1G:79:G:N3	2.26	0.50
20:BA:86:ARG:CZ	20:BA:86:ARG:HB2	2.41	0.50
2:12:12:GLU:HB2	2:12:16:HIS:CG	2.46	0.50
55:Q8:46:ARG:NH2	55:Q8:48:PHE:HA	2.25	0.50
24:3K:13:C:O2'	24:3K:14:A:H5'	2.10	0.50
13:4A:70:LEU:O	13:4A:74:VAL:HG23	2.11	0.50
2:12:163:PHE:CD1	2:12:185:ILE:HG13	2.44	0.50
5:42:101:ILE:O	5:42:120:THR:OG1	2.19	0.50
8:72:120:THR:OG1	8:72:121:ASP:N	2.44	0.50
7:6E:15:ASP:OD1	7:6E:16:LEU:N	2.44	0.50
26:1H:2698:U:H2'	26:1H:2699:C:C6	2.46	0.50
43:A5:73:ALA:O	43:A5:106:ILE:HG12	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:1E:111:ARG:NH1	2:1E:111:ARG:HG2	2.27	0.50
40:75:107:ASP:OD1	40:75:109:GLU:HB2	2.11	0.50
26:1H:2347:C:P	53:O8:39:TYR:OH	2.70	0.50
32:51:12:PRO:HG2	32:51:13:LYS:HE2	1.93	0.50
4:3E:175:SER:O	4:3E:183:GLY:HA2	2.12	0.50
29:29:120:TRP:CD1	29:29:155:LYS:HB3	2.46	0.50
26:1H:6:A:O2'	34:58:129:PRO:HB3	2.12	0.50
50:H5:7:LYS:HG3	50:H5:34:GLU:HG3	1.92	0.50
26:14:1708:C:O2'	26:14:1709:U:H5'	2.11	0.50
9:8E:4:TYR:CZ	9:8E:88:TYR:HB2	2.46	0.50
1:1G:298:A:H5''	1:1G:299:G:OP2	2.11	0.50
20:BA:50:GLU:HA	20:BA:100:ILE:HG21	1.93	0.50
31:41:118:ARG:O	31:41:181:ARG:HG3	2.12	0.50
1:1G:967:C:H3'	1:1G:968:A:H2'	1.94	0.50
18:9I:52:PRO:O	18:9I:56:THR:HG23	2.10	0.50
25:4K:13:A:O2'	25:4K:14:A:OP1	2.30	0.50
33:69:90:GLY:O	33:69:121:LYS:HD2	2.11	0.50
26:14:30:G:H2'	26:14:31:C:C6	2.46	0.50
43:A5:88:ARG:HD3	43:A5:94:ASP:OD2	2.11	0.50
26:14:832:G:H5'	36:35:45:LEU:HD12	1.94	0.50
24:3K:48:C:C5	24:3K:59:U:H1'	2.47	0.50
44:F8:3:THR:CB	44:F8:4:ALA:HA	2.41	0.50
42:95:76:LYS:HD2	42:95:80:GLN:O	2.10	0.50
34:15:35:ARG:HB2	34:15:42:TRP:CZ3	2.47	0.50
28:19:32:SER:O	28:19:33:LEU:HB2	2.11	0.50
28:19:33:LEU:HD21	28:19:103:ARG:HA	1.93	0.50
26:14:882:G:N2	26:14:894:C:H42	2.08	0.50
53:O8:28:ARG:CZ	53:O8:30:THR:HG23	2.41	0.50
26:14:2646:C:H2'	26:14:2647:U:O4'	2.11	0.50
1:1G:373:A:N3	1:1G:374:A:C8	2.80	0.50
2:12:82:ARG:HG3	2:12:92:TYR:OH	2.10	0.50
36:35:78:PRO:HA	36:35:110:TYR:CD2	2.46	0.50
26:14:2184:G:H2'	26:14:2185:C:C6	2.45	0.50
20:BA:64:ASP:OD1	20:BA:81:LYS:HD2	2.11	0.50
54:L5:12:ARG:NH2	54:L5:44:PRO:HB3	2.27	0.50
18:9A:70:ILE:HG23	18:9A:79:LEU:HD12	1.92	0.50
26:1H:1718:G:C2	26:1H:1725:G:C8	2.99	0.50
26:1H:390:A:C6	36:78:71:VAL:HG21	2.47	0.50
26:14:1507:A:C5	26:14:1508:A:H1'	2.46	0.50
26:1H:1216:G:OP2	41:C8:12:ARG:NH2	2.42	0.50
26:1H:1424:G:H2'	26:1H:1425:G:O4'	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
43:A5:79:GLY:HA3	43:A5:100:THR:HG22	1.94	0.50
26:1H:1187:G:HO2'	26:1H:1188:U:H6	1.57	0.50
55:M5:40:GLU:H	55:M5:43:GLN:HG3	1.77	0.50
50:L8:7:LYS:HA	50:L8:33:GLN:O	2.12	0.50
1:1G:1352:C:N4	1:1G:1370:G:H1	2.08	0.50
26:1H:1534:G:H21	26:1H:1535:U:H5''	1.76	0.50
49:K8:47:ASN:C	49:K8:49:LYS:H	2.14	0.50
26:14:819:A:C4	26:14:1189:A:C2	2.99	0.50
5:42:37:ARG:HG2	5:42:112:LEU:HA	1.93	0.50
1:13:1260:C:H6	1:13:1260:C:H3'	1.76	0.50
26:14:1011:G:C2	26:14:1151:G:C2	3.00	0.50
8:7E:109:ILE:HD11	8:7E:120:THR:CG2	2.42	0.50
35:68:120:GLU:HB2	40:B8:68:TYR:CE2	2.46	0.50
40:75:55:ASN:N	40:75:59:THR:HG22	2.26	0.50
53:K5:28:ARG:HB3	53:K5:32:ASN:H	1.77	0.50
26:1H:518:G:H2'	26:1H:519:U:H6	1.76	0.50
26:14:2693:A:H2'	26:14:2694:G:C8	2.47	0.50
24:3L:54:U:H3	24:3L:58:A:H62	1.59	0.50
5:42:79:GLU:OE1	8:72:104:ARG:HA	2.11	0.50
1:1G:1387:G:H2'	1:1G:1388:C:H6	1.76	0.50
5:42:24:ARG:HB3	5:42:24:ARG:CZ	2.41	0.50
23:2L:54:G:H2'	23:2L:55:5MU:H6	1.76	0.50
26:14:1298:C:H5''	26:14:1299:G:OP2	2.11	0.50
40:75:29:ARG:HD3	40:75:44:ASP:OD2	2.11	0.50
1:1G:748:C:H4'	1:1G:749:C:O5'	2.12	0.50
29:21:3:GLY:HA3	29:21:81:ILE:HG21	1.94	0.50
26:1H:50:U:H3'	26:1H:51:G:H5'	1.93	0.50
26:1H:2246:G:H2'	26:1H:2247:A:C8	2.46	0.50
26:1H:320:A:H2'	30:31:136:THR:HG21	1.94	0.50
28:11:52:ARG:HB2	28:11:53:PHE:CD2	2.46	0.50
10:1I:96:ILE:H	10:1I:96:ILE:HD13	1.77	0.50
1:13:807:A:H2'	1:13:808:C:C6	2.46	0.50
26:1H:336:C:OP1	45:G8:83:THR:HG23	2.11	0.50
26:14:1044:G:O2'	26:14:1045:A:H5''	2.11	0.50
46:H8:124:ILE:HD12	46:H8:125:LEU:H	1.76	0.50
44:F8:3:THR:HB	44:F8:6:ASP:HB2	1.93	0.50
26:14:2074:U:H2'	26:14:2075:U:C6	2.47	0.50
26:1H:140:A:C8	26:1H:1408:C:O2'	2.63	0.50
26:14:273(C):C:H5'	26:14:273(D):C:OP2	2.11	0.50
41:C8:92:ARG:NH1	42:D8:11:GLN:O	2.45	0.50
26:1H:1062:G:N2	26:1H:1076:C:O2	2.44	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:94:G:N3	49:G5:47:ASN:ND2	2.59	0.50
46:D5:69:THR:HB	46:D5:88:PHE:HB3	1.93	0.50
26:14:1532:C:H42	26:14:1539:G:H1	1.59	0.50
26:1H:270(E):G:C6	26:1H:270(F):U:C4	2.99	0.50
37:45:25:ASP:HB3	37:45:102:VAL:HG23	1.93	0.50
26:14:654(R):C:N4	26:14:654(S):G:O6	2.44	0.50
2:1E:9:GLU:HA	2:1E:12:GLU:HB2	1.93	0.50
1:1G:1329:A:H5'	13:4A:29:ARG:NE	2.26	0.50
1:1G:690:G:H22	11:2A:55:LYS:HE2	1.77	0.50
1:1G:631:G:H3'	1:1G:632:A:H8	1.76	0.50
26:1H:586:A:OP1	58:1H:3827:HOH:O	2.20	0.50
1:13:627:G:H2'	1:13:628:G:H8	1.76	0.50
45:G8:38:ILE:HD11	45:G8:64:GLU:HG3	1.93	0.50
44:F8:11:PRO:HD3	49:K8:37:PHE:CD2	2.47	0.50
26:1H:7:G:H2'	26:1H:8:A:O4'	2.11	0.50
37:45:34:LEU:HD12	37:45:130:LYS:O	2.12	0.50
8:7E:110:ALA:HB3	8:7E:121:ASP:HB3	1.92	0.50
46:H8:48:PHE:HE1	46:H8:71:VAL:HG11	1.76	0.50
26:1H:1843:C:H5'	28:11:253:GLN:OE1	2.11	0.50
26:1H:1013:C:O2'	26:1H:1014:U:H5'	2.12	0.50
9:8E:86:VAL:O	9:8E:90:PRO:HB3	2.11	0.50
40:B8:84:GLN:HG2	40:B8:85:LYS:HG2	1.92	0.50
6:5E:69:GLU:HG2	6:5E:70:ASP:H	1.77	0.50
26:14:2540:C:O2'	26:14:2740:A:N3	2.43	0.50
26:1H:259:G:N2	26:1H:621:A:H8	2.06	0.50
24:3K:58:A:HO2'	24:3K:59:U:P	2.35	0.50
30:39:112:MET:O	30:39:115:ALA:HB3	2.12	0.50
26:14:2134:A:N7	26:14:2157:G:H1'	2.26	0.50
26:14:1434:A:H2'	26:14:1435:G:C8	2.47	0.50
1:1G:742:G:H5''	15:6A:58:MET:HE1	1.93	0.50
51:M8:23:GLU:OE1	51:M8:24:THR:N	2.45	0.50
22:1K:27:G:H2'	22:1K:28:G:C8	2.46	0.50
40:B8:111:ARG:H	40:B8:111:ARG:CD	2.24	0.50
29:29:128:SER:OG	29:29:129:HIS:N	2.43	0.50
28:11:271:ILE:O	28:11:272:ALA:HB2	2.11	0.50
2:1E:126:GLU:HA	2:1E:129:GLU:CG	2.42	0.50
26:14:2675:A:H4'	35:25:29:ASN:ND2	2.26	0.50
3:2E:130:VAL:O	3:2E:134:ILE:HG12	2.11	0.50
26:14:2302:G:O2'	31:49:126:ASP:HB3	2.10	0.50
1:13:984:C:N4	1:13:1221:G:H1	2.09	0.50
23:2K:33:OMC:HM22	23:2K:34:U:H5'	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:194:G:H2'	26:14:195:A:O4'	2.12	0.50
35:25:24:VAL:HG23	35:25:33:ALA:HB2	1.94	0.50
26:14:1680:U:N3	26:14:1764:G:OP2	2.31	0.50
26:1H:2864:G:OP1	40:B8:119:LYS:HD2	2.12	0.50
6:52:72:VAL:O	6:52:75:LEU:HB3	2.12	0.50
34:15:13:TRP:O	34:15:135:PRO:HD2	2.12	0.50
26:1H:1864:U:OP1	26:1H:2410:G:O2'	2.23	0.50
17:8I:48:GLU:O	17:8I:50:LYS:HG2	2.12	0.50
26:14:2292:C:OP1	39:65:17:ARG:NH2	2.43	0.50
1:13:371:G:O2'	1:13:373:A:N7	2.43	0.50
24:3K:75:C:O2'	24:3K:76:A:H2	1.94	0.50
1:13:1399:C:C2	1:13:1502:A:N6	2.80	0.50
26:14:1111:A:O3'	32:59:3:ARG:HB3	2.11	0.50
26:1H:2583:G:OP1	58:1H:4456:HOH:O	2.20	0.50
29:29:81:ILE:HG21	29:29:84:PHE:HD2	1.77	0.50
55:M5:22:VAL:HB	55:M5:53:PRO:HG3	1.94	0.50
23:2K:48:U:O2'	23:2K:49:C:OP2	2.26	0.50
46:H8:116:VAL:O	46:H8:174:VAL:HA	2.11	0.50
1:1G:538:G:H2'	1:1G:539:A:C8	2.46	0.50
26:1H:198:C:H4'	26:1H:2243:U:O2'	2.12	0.50
37:88:134:ARG:NH2	46:H8:122:ARG:HD2	2.27	0.50
1:13:444:C:H2'	1:13:445:G:H8	1.77	0.50
37:45:31:ASP:O	37:45:134:ARG:HB3	2.11	0.50
26:1H:1296:G:O2'	26:1H:1297:C:H5'	2.11	0.50
4:3E:108:LEU:HD23	4:3E:110:PHE:HE1	1.77	0.50
5:4E:15:ARG:HH11	25:4K:25:A:H3'	1.75	0.50
42:D8:24:LYS:HA	42:D8:92:THR:HG23	1.94	0.50
29:29:120:TRP:CD2	29:29:155:LYS:HD3	2.47	0.50
1:1G:1386:G:C2	1:1G:1387:G:N7	2.80	0.50
26:14:389:G:H22	36:35:72:PRO:CD	2.25	0.50
1:13:300:A:H1'	1:13:565:U:O2	2.11	0.50
45:G8:78:ALA:HB3	45:G8:79:CYS:SG	2.52	0.50
16:7A:19:ILE:HB	16:7A:36:ILE:O	2.11	0.50
26:14:1779:U:H5''	26:14:1780:A:H5'	1.94	0.50
18:9I:32:ARG:NH1	18:9I:65:ILE:HD13	2.27	0.50
26:14:2567:G:H2'	26:14:2568:C:C6	2.47	0.50
24:1L:2:C:N3	24:1L:72:C:N4	2.60	0.50
26:1H:1500:G:O2'	28:11:100:GLY:O	2.26	0.50
44:B5:56:THR:HB	44:B5:77:LYS:HE2	1.94	0.50
32:59:163:TYR:CZ	32:59:169:VAL:HG21	2.47	0.50
24:3K:46:G:O2'	24:3K:48:C:O2	2.30	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
29:29:70:ALA:HB1	29:29:72:VAL:HG23	1.94	0.50
26:14:900:A:H3'	26:14:901:A:C8	2.45	0.50
1:1G:464:G:C6	1:1G:466:C:H5'	2.47	0.50
26:1H:963:U:H5''	58:1H:3796:HOH:O	2.11	0.50
26:14:2394:C:OP1	36:35:63:PRO:HG2	2.11	0.50
29:21:48:GLN:OE1	29:21:77:ILE:HG21	2.11	0.50
30:39:128:ALA:O	30:39:142:TRP:NE1	2.45	0.50
29:21:19:ARG:HA	35:68:73:ASP:HA	1.93	0.50
32:51:154:PRO:HB3	32:51:163:TYR:CZ	2.47	0.50
5:4E:12:LEU:HB3	5:4E:31:LEU:HB2	1.93	0.50
5:42:122:GLU:HG2	5:42:131:ILE:HD12	1.94	0.50
13:4A:78:ILE:HD13	13:4A:92:HIS:CE1	2.47	0.50
26:14:2823:A:OP1	29:29:159:HIS:NE2	2.42	0.50
51:I5:56:VAL:HG22	51:I5:57:GLU:HG3	1.94	0.50
26:1H:280:C:N3	26:1H:361:G:C2	2.80	0.50
24:3K:71:G:O2'	26:1H:1851:U:O2'	2.26	0.50
30:31:178:PRO:HB2	30:31:201:VAL:HG21	1.94	0.50
26:1H:1288:U:H4'	26:1H:1289:C:OP2	2.11	0.50
1:1G:1018:C:H2'	1:1G:1019:C:O4'	2.12	0.50
1:1G:685:G:C2	1:1G:686:U:C4	2.99	0.50
55:M5:30:ARG:O	55:M5:32:LEU:N	2.44	0.50
20:BA:79:ARG:HE	20:BA:83:ARG:NH1	2.10	0.50
26:14:354:G:H2'	26:14:355:G:H8	1.76	0.50
26:14:1630:G:N2	26:14:1636:C:O2	2.33	0.50
9:82:49:PRO:HB2	9:82:85:LEU:HD21	1.94	0.50
26:14:531:C:C5	26:14:2035:G:C2	3.00	0.49
26:1H:2275:C:H5'	26:1H:2275:C:C6	2.47	0.49
5:42:102:ALA:HB3	5:42:107:ARG:HB2	1.94	0.49
1:13:963:G:H21	10:1I:55:LYS:HZ1	1.60	0.49
10:1I:49:VAL:HG23	14:5I:41:ARG:HB2	1.94	0.49
1:1G:672:U:H2'	1:1G:673:G:C8	2.47	0.49
1:1G:427:U:H3'	1:1G:428:G:H2'	1.93	0.49
1:13:522:C:H2'	1:13:523:A:O4'	2.12	0.49
51:M8:21:VAL:O	51:M8:24:THR:HG23	2.11	0.49
1:1G:848:C:H2'	1:1G:849:C:C6	2.47	0.49
13:4A:3:ARG:HB2	51:I5:34:GLU:CG	2.40	0.49
26:1H:229:A:OP2	36:78:150:ALA:HB1	2.11	0.49
26:14:1340:U:H4'	26:14:1341:U:OP2	2.11	0.49
5:42:7:GLU:OE1	5:42:37:ARG:NH2	2.45	0.49
1:13:624:C:H4'	16:7I:11:SER:N	2.26	0.49
26:14:2151:G:H2'	26:14:2152:G:C8	2.47	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:13:592:G:H1	1:13:647:C:H42	1.58	0.49
33:61:69:LYS:HA	33:61:136:VAL:HB	1.94	0.49
26:14:1027:A:H5'	27:1J:88:C:N4	2.27	0.49
26:14:321:G:OP1	30:39:135:LYS:NZ	2.31	0.49
26:14:2542:A:H4'	26:14:2542:A:OP1	2.11	0.49
1:13:1192:C:OP2	3:2E:4:LYS:NZ	2.44	0.49
1:13:67:C:H2'	1:13:68:G:H8	1.75	0.49
26:14:2301:C:H2'	26:14:2302:G:H8	1.77	0.49
4:3E:162:LEU:O	4:3E:165:MET:HB3	2.12	0.49
15:6I:38:ARG:HH11	15:6I:38:ARG:HG2	1.77	0.49
37:45:34:LEU:HD11	37:45:129:THR:HB	1.94	0.49
41:85:8:VAL:O	41:85:12:ARG:HG2	2.12	0.49
1:1G:577:G:C8	1:1G:816:A:C6	3.00	0.49
1:1G:1194:U:H2'	1:1G:1195:C:C6	2.47	0.49
1:1G:1139:G:H1	1:1G:1143:G:H22	1.59	0.49
26:14:2412:A:H2'	26:14:2413:G:O4'	2.12	0.49
27:16:21:G:H1	27:16:62:C:H42	1.58	0.49
26:1H:492:A:H2'	26:1H:493:G:O4'	2.12	0.49
7:6E:150:ALA:HB2	11:2I:50:TYR:OH	2.12	0.49
26:14:988:A:N6	50:H5:13:ILE:HG21	2.27	0.49
30:39:160:ASN:HB3	30:39:163:VAL:HB	1.94	0.49
16:7I:67:THR:H	16:7I:70:ALA:HB3	1.77	0.49
37:88:112:GLU:CD	37:88:112:GLU:H	2.15	0.49
26:1H:997:G:OP1	41:C8:93:LYS:HD2	2.11	0.49
55:Q8:50:LEU:O	55:Q8:51:ALA:HB3	2.12	0.49
24:3K:45:U:H1'	24:3K:46:G:H5''	1.94	0.49
27:1J:72:G:O2'	27:1J:104:A:N6	2.45	0.49
8:7E:6:ILE:HB	8:7E:85:ARG:HH12	1.76	0.49
26:1H:2807:G:H3'	26:1H:2808:U:H5''	1.94	0.49
55:M5:34:TRP:CG	55:M5:35:GLN:N	2.81	0.49
26:1H:1580:A:OP2	26:1H:1580:A:H8	1.96	0.49
26:14:2262:U:O2'	26:14:2263:C:H5'	2.13	0.49
1:1G:1015:A:O5'	1:1G:1015:A:H8	1.95	0.49
28:11:183:ARG:NH1	28:11:269:PHE:HB2	2.27	0.49
1:13:342:C:N4	1:13:347:G:H1	2.10	0.49
26:1H:270(L):U:N3	33:61:50:ARG:HG2	2.26	0.49
26:14:654(B):C:H2'	26:14:654(C):G:C8	2.46	0.49
26:1H:582:G:H2'	26:1H:583:G:C8	2.48	0.49
37:45:75:THR:HG23	37:45:89:ASN:HB3	1.94	0.49
1:13:444:C:H2'	1:13:445:G:C8	2.47	0.49
28:19:37:LEU:HD13	28:19:39:LYS:HB2	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
20:BA:26:ASN:O	20:BA:30:LYS:HB2	2.12	0.49
11:2I:45:GLY:O	11:2I:50:TYR:HB2	2.12	0.49
26:14:1657:C:H2'	26:14:1658:C:C6	2.46	0.49
1:1G:736:C:H2'	1:1G:737:A:C8	2.46	0.49
26:1H:2693:A:H2'	26:1H:2694:G:H8	1.76	0.49
45:C5:84:ARG:NH2	45:C5:94:LYS:HE2	2.27	0.49
26:1H:128:C:H2'	26:1H:129:C:H6	1.77	0.49
1:1G:1147:C:O2	9:82:16:ARG:NE	2.46	0.49
8:72:30:ARG:O	8:72:34:GLU:HG2	2.13	0.49
46:D5:60:GLU:HB2	46:D5:66:SER:OG	2.12	0.49
48:J8:92:LYS:HG3	48:J8:95:LEU:HD12	1.94	0.49
36:35:127:ALA:O	36:35:147:LEU:N	2.45	0.49
36:35:146:VAL:HG13	36:35:147:LEU:HD22	1.93	0.49
26:1H:598:G:H1'	36:78:12:ALA:HB2	1.94	0.49
16:7I:57:ARG:NH2	16:7I:79:VAL:O	2.44	0.49
1:13:1128:C:O2'	1:13:1129:C:H5''	2.12	0.49
24:1L:18:G:O2'	24:1L:19:G:OP1	2.30	0.49
1:13:129(A):G:N2	1:13:188:U:HO2'	2.10	0.49
26:14:633:A:O2'	26:14:2404:C:OP1	2.28	0.49
26:1H:2635:C:H5''	29:21:78:LEU:HA	1.94	0.49
17:8I:29:HIS:CD2	17:8I:30:PRO:HD2	2.47	0.49
26:14:975:G:C2	26:14:990:A:C8	2.99	0.49
19:AI:19:VAL:HG13	19:AI:47:HIS:CD2	2.45	0.49
26:14:2346:A:H5''	26:14:2383:G:O4'	2.12	0.49
1:13:792:A:H4'	1:13:793:U:O5'	2.12	0.49
1:1G:601:C:H2'	1:1G:602:A:C8	2.48	0.49
47:I8:38:VAL:HG12	47:I8:40:GLN:HG2	1.95	0.49
29:29:117:MET:HA	29:29:122:PHE:N	2.28	0.49
26:1H:821:A:H2'	26:1H:946:G:H5''	1.94	0.49
4:32:178:VAL:C	4:32:180:GLY:H	2.16	0.49
26:1H:172:C:H2'	26:1H:173:G:C8	2.46	0.49
26:14:1005:C:O2'	34:15:28:THR:HG21	2.12	0.49
1:13:1418:A:C2	1:13:1483:A:C2	3.01	0.49
46:D5:40:ASP:HB3	46:D5:43:GLU:HB2	1.94	0.49
2:1E:88:ALA:HB2	2:1E:219:VAL:HG13	1.94	0.49
39:A8:67:ARG:HB2	39:A8:67:ARG:NH1	2.27	0.49
30:39:21:ALA:C	30:39:23:ASP:H	2.16	0.49
26:14:305:U:H2'	26:14:306:U:C6	2.46	0.49
26:14:2239:G:P	58:14:3510:HOH:O	2.69	0.49
26:14:1693:U:O2'	28:19:14:ARG:NH2	2.45	0.49
28:11:172:TYR:CD2	28:11:186:HIS:HA	2.46	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1G:410:G:N1	1:1G:429:U:O2	2.45	0.49
1:13:994:A:N3	1:13:994:A:H2'	2.27	0.49
26:14:1716:U:H2'	26:14:1717:G:H8	1.76	0.49
1:13:1363:A:H1'	1:13:1365:G:N7	2.27	0.49
26:14:1083:U:H1'	26:14:1086:A:N6	2.28	0.49
28:11:67:PHE:HB3	28:11:153:ALA:H	1.77	0.49
28:11:65:ILE:HD11	28:11:67:PHE:CE1	2.48	0.49
32:59:82:GLY:HA3	32:59:135:GLY:O	2.11	0.49
1:1G:620:C:C2	4:32:135:LEU:HG	2.48	0.49
1:13:43:C:HO2'	1:13:623:C:HO2'	1.60	0.49
1:1G:1443:G:O2'	40:75:122:ASP:OD2	2.30	0.49
26:14:71:A:H4'	26:14:72:U:H5''	1.94	0.49
26:14:1021:A:H3'	26:14:1021:A:C8	2.46	0.49
26:14:1754:C:H2'	26:14:1755:A:C8	2.47	0.49
27:1J:38:C:O4'	39:65:95:HIS:NE2	2.45	0.49
26:14:2352:A:H2	47:E5:33:ALA:HB1	1.72	0.49
50:L8:50:VAL:O	50:L8:54:VAL:HG12	2.13	0.49
32:51:4:ILE:C	32:51:6:ARG:H	2.14	0.49
1:13:79:G:N2	1:13:89:U:H2'	2.26	0.49
26:14:2262:U:OP2	47:E5:19:LYS:HE2	2.12	0.49
28:19:255:LYS:H	28:19:255:LYS:HE3	1.77	0.49
26:14:2296:U:OP2	39:65:9:ARG:NH1	2.44	0.49
1:1G:1299:A:C6	1:1G:1301:U:C2	3.00	0.49
3:22:15:THR:HG21	3:22:181:ASN:HA	1.94	0.49
1:1G:1086:U:H3	1:1G:1099:G:H22	1.60	0.49
26:14:918:A:O2'	27:1J:96:G:N2	2.45	0.49
1:13:209:U:H5'	1:13:210:U:OP2	2.12	0.49
24:3K:24:G:H2'	24:3K:25:C:C6	2.46	0.49
24:3K:24:G:H2'	24:3K:25:C:H6	1.77	0.49
27:1J:9:G:H5'	39:65:25:ARG:HH12	1.77	0.49
28:11:89:SER:HB2	28:11:159:ALA:HB2	1.94	0.49
4:3E:155:LEU:O	4:3E:158:ILE:N	2.45	0.49
26:14:2889:C:H3'	26:14:2891:G:H8	1.77	0.49
1:13:1352:C:H2'	1:13:1353:G:C8	2.47	0.49
43:E8:57:ASN:O	43:E8:61:ASN:HB2	2.13	0.49
26:1H:2010:G:N7	58:1H:4106:HOH:O	2.35	0.49
26:1H:1641:A:H2'	26:1H:1642:G:O4'	2.12	0.49
26:1H:2209:C:O2	26:1H:2216:G:C2	2.66	0.49
26:14:2241:A:H2'	26:14:2242:G:C8	2.47	0.49
10:1I:5:ARG:HB2	10:1I:73:ASP:OD1	2.13	0.49
1:1G:129:U:O4	58:1G:1763:HOH:O	2.18	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
31:49:60:LEU:O	31:49:64:THR:HG22	2.11	0.49
1:1G:111:G:H8	1:1G:111:G:O5'	1.96	0.49
26:1H:1301:A:N3	26:1H:1301:A:H2'	2.28	0.49
12:3I:9:GLN:O	12:3I:13:LYS:HG2	2.12	0.49
26:14:1542:G:O5'	26:14:1543:A:H5''	2.11	0.49
1:13:284:G:H2'	1:13:285:G:C8	2.48	0.49
24:1L:68:C:H2'	24:1L:69:G:C8	2.46	0.49
1:13:1187:G:O5'	9:8E:113:LYS:NZ	2.45	0.49
26:1H:2127:G:N2	26:1H:2162:G:H1'	2.18	0.49
4:3E:11:LEU:HD22	4:3E:66:ARG:HG2	1.94	0.49
27:16:12:C:C2	47:I8:74:ARG:NH1	2.81	0.49
32:59:66:GLY:O	32:59:70:THR:OG1	2.29	0.49
33:69:73:GLU:HG3	33:69:136:VAL:HG23	1.94	0.49
26:14:2376:A:H2	39:65:112:PHE:HB2	1.78	0.49
1:13:258:G:H2'	1:13:259:G:H8	1.77	0.49
26:14:882:G:H1	26:14:894:C:H42	1.60	0.49
1:1G:644:G:H4'	8:72:92:ARG:NH1	2.27	0.49
29:29:24:THR:HG21	29:29:188:VAL:HG12	1.93	0.49
26:14:2320:A:C6	26:14:2333:A:C8	3.00	0.49
26:1H:270(L):U:C2	33:61:50:ARG:HG2	2.47	0.49
9:8E:26:VAL:HB	9:8E:33:PHE:HB2	1.95	0.49
1:13:64:G:H4'	1:13:65:U:H5'	1.94	0.49
26:1H:2531:A:H5'	32:51:157:TYR:CZ	2.47	0.49
26:1H:2111:C:O2'	26:1H:2119:A:OP1	2.29	0.49
1:13:458:C:N4	1:13:475:G:H1	2.10	0.49
48:F5:52:ARG:CZ	48:F5:56:GLN:HA	2.42	0.49
1:13:1489:G:H2'	1:13:1490:C:O4'	2.13	0.49
9:82:53:VAL:HG23	9:82:55:ALA:H	1.76	0.49
28:11:76:PRO:HG2	28:11:98:VAL:HG11	1.94	0.49
6:5E:3:ARG:HB3	6:5E:93:SER:HB2	1.94	0.49
23:2L:62:C:H2'	23:2L:63:C:H6	1.78	0.49
2:12:71:VAL:HB	2:12:170:GLU:HG2	1.95	0.49
26:14:110:G:C2	26:14:111:A:C8	3.01	0.49
26:1H:2309:A:C5	26:1H:2310:A:H8	2.30	0.49
3:22:120:VAL:HG21	3:22:137:ALA:HB2	1.94	0.49
1:1G:692:U:H5	11:2A:26:ASN:OD1	1.96	0.49
7:62:149:ARG:HD3	11:2A:59:TYR:CE1	2.47	0.49
39:65:106:ARG:O	39:65:106:ARG:HD2	2.12	0.49
1:13:45:U:O5'	1:13:45:U:H6	1.95	0.49
43:A5:71:VAL:HA	43:A5:107:LEU:HD12	1.95	0.49
1:1G:447:G:O2'	1:1G:487:A:N6	2.45	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:1H:1621:U:H5''	26:1H:1622:G:OP1	2.13	0.49
26:1H:1188:U:C4'	42:D8:79:VAL:HG22	2.42	0.49
7:6E:113:GLU:HB2	7:6E:118:VAL:HG13	1.94	0.49
26:1H:2053:G:OP1	29:21:144:ARG:HD3	2.12	0.49
26:1H:445:C:O2'	26:1H:446:G:H5'	2.13	0.49
47:E5:49:LYS:HE3	47:E5:82:ARG:CZ	2.43	0.49
27:1J:38:C:N4	27:1J:44:G:H1	2.05	0.49
26:14:921:G:C6	26:14:922:U:C4	3.01	0.49
33:61:40:THR:O	33:61:44:LEU:HB2	2.13	0.49
26:14:2690:C:OP1	38:55:17:ARG:NH1	2.44	0.49
26:1H:991:C:H42	26:1H:1163:G:H1	1.61	0.49
1:1G:728:A:C2	1:1G:729:A:C5	3.01	0.49
26:1H:2059:A:H5''	26:1H:2060:A:OP2	2.12	0.49
13:4I:7:VAL:HB	31:41:115:ARG:NH2	2.28	0.49
1:13:1309:G:OP1	13:4I:88:ARG:NH1	2.46	0.49
42:95:48:GLY:HA3	42:95:52:VAL:N	2.28	0.49
29:29:5:LEU:HD23	29:29:51:PHE:HB2	1.95	0.49
26:14:289:A:H3'	26:14:290:G:C8	2.46	0.49
1:1G:559:A:H4'	1:1G:560:U:H5''	1.93	0.49
26:14:1190:G:O2'	26:14:1191:G:H5'	2.12	0.49
45:G8:28:LYS:NZ	45:G8:64:GLU:OE2	2.28	0.49
26:14:2889:C:H3'	26:14:2891:G:C8	2.47	0.49
32:51:92:ILE:CD1	32:51:93:GLY:H	2.25	0.49
24:3K:71:G:HO2'	26:1H:1851:U:HO2'	1.55	0.49
26:14:2116:G:H2'	26:14:2117:A:C4	2.48	0.49
26:14:755:C:H2'	26:14:756:C:C6	2.48	0.49
26:14:1480:G:C6	26:14:1482:U:C4	3.00	0.49
2:12:222:ILE:O	2:12:226:ARG:HB2	2.13	0.49
26:14:857:C:H2'	26:14:858:U:H6	1.77	0.49
26:1H:2688:U:H5	26:1H:2720:U:OP2	1.96	0.49
13:4I:91:ARG:HB2	13:4I:98:VAL:HG23	1.94	0.49
26:1H:2404:C:O3'	36:78:77:ARG:NH2	2.45	0.49
4:32:150:GLU:C	4:32:152:SER:H	2.16	0.49
27:1J:53:A:H2'	27:1J:54:G:O4'	2.12	0.49
55:Q8:8:LYS:HD2	55:Q8:8:LYS:N	2.27	0.49
35:25:98:VAL:HG12	35:25:117:LEU:HB3	1.93	0.49
35:25:98:VAL:CG1	35:25:117:LEU:HB3	2.42	0.49
1:1G:440:A:H3'	1:1G:442:C:C6	2.47	0.49
1:1G:1266:G:N2	1:1G:1270:C:N3	2.61	0.49
1:13:1392:G:H21	1:13:1502:A:H8	1.61	0.49
26:1H:1729:A:C6	26:1H:1731:G:C2	3.00	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:12:67:THR:H	2:12:160:ASP:HB2	1.78	0.49
1:13:737:A:O2'	1:13:738:C:H5'	2.13	0.49
26:1H:2056:G:H2'	26:1H:2056:G:N3	2.28	0.49
26:14:2638:G:OP1	29:29:82:ARG:NH2	2.45	0.49
1:1G:1067:A:H4'	1:1G:1068:G:O5'	2.13	0.49
33:61:40:THR:HG22	33:61:41:GLU:H	1.77	0.49
41:85:100:VAL:O	41:85:101:ARG:HG2	2.12	0.49
23:2K:54:G:C5	23:2K:55:5MU:H72	2.47	0.49
40:75:8:LYS:HZ2	40:75:8:LYS:CB	2.24	0.49
29:29:111:ARG:HB2	29:29:160:TYR:O	2.12	0.49
15:6A:87:ILE:HG22	15:6A:88:ARG:N	2.25	0.49
13:4A:84:ILE:O	13:4A:86:CYS:N	2.46	0.49
39:A8:83:LYS:HG3	39:A8:110:LEU:CD1	2.43	0.49
30:39:129:PHE:HA	30:39:142:TRP:CD1	2.48	0.49
39:A8:59:LYS:HG2	39:A8:60:GLY:N	2.28	0.49
26:1H:1025:G:C4	26:1H:1135:C:H1'	2.47	0.49
2:1E:69:LEU:HB3	2:1E:162:ILE:HG22	1.95	0.49
26:14:4:C:N4	26:14:2899:G:H1	2.11	0.49
10:1A:4:ILE:HG13	10:1A:77:PRO:HB3	1.95	0.49
29:21:51:PHE:O	29:21:74:PRO:HB2	2.11	0.49
1:1G:485:G:O2'	1:1G:486:U:H6	1.96	0.49
26:14:1171:G:O2'	26:14:1173:G:N3	2.46	0.49
30:39:83:PHE:O	30:39:84:VAL:HB	2.13	0.49
47:I8:23:VAL:HG13	47:I8:38:VAL:HG23	1.95	0.49
26:1H:930:U:H4'	26:1H:931:G:O5'	2.12	0.49
21:1B:2:GLY:O	21:1B:5:ASP:N	2.34	0.49
44:F8:11:PRO:HB3	44:F8:92:LEU:HD21	1.95	0.49
1:13:163:C:H2'	1:13:164:U:C6	2.48	0.49
1:13:1503:A:O2'	25:4K:13:A:N6	2.45	0.49
1:1G:403:C:N4	1:1G:547:A:H5'	2.28	0.49
49:G5:43:GLN:HB2	49:G5:45:SER:H	1.76	0.49
18:9A:19:LYS:HG3	18:9A:20:ALA:H	1.78	0.49
32:59:103:LEU:HD23	32:59:103:LEU:H	1.78	0.49
18:9I:38:GLU:HA	18:9I:41:LYS:NZ	2.28	0.49
26:1H:687:C:OP2	58:1H:4021:HOH:O	2.20	0.49
26:14:984:A:H5''	26:14:985:C:H5	1.77	0.49
37:45:59:ARG:O	37:45:60:ARG:HG3	2.12	0.49
1:13:192:U:H2'	1:13:193:C:C6	2.48	0.49
1:1G:1162:C:H42	1:1G:1174:G:H1	1.61	0.49
1:1G:1126:U:H4'	1:1G:1127:G:H8	1.77	0.49
26:1H:2395:C:H5''	26:1H:2396:G:OP2	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
39:A8:35:ILE:HD11	39:A8:101:LEU:HD23	1.93	0.49
39:A8:36:TYR:N	39:A8:36:TYR:CD1	2.81	0.49
1:13:111:G:H5''	16:7I:27:LYS:HG2	1.95	0.49
1:1G:1036:G:H5'	1:1G:1037:C:OP2	2.12	0.49
3:22:6:HIS:HB3	14:5A:49:HIS:HD2	1.76	0.49
26:1H:265:A:H1'	26:1H:266:G:O4'	2.12	0.49
49:K8:59:ARG:O	49:K8:62:THR:HG23	2.13	0.49
13:4A:3:ARG:HG2	13:4A:8:GLU:O	2.13	0.49
8:72:92:ARG:HB3	8:72:94:TYR:CE2	2.48	0.49
38:98:72:ASP:O	38:98:76:VAL:HG23	2.13	0.49
26:1H:2262:U:OP1	26:1H:2387:U:O2'	2.19	0.49
26:1H:2729:G:N3	29:21:187:ALA:HB2	2.27	0.49
26:14:1011:G:H2'	26:14:1013:C:O4'	2.13	0.49
26:14:868:U:H2'	26:14:869:G:C8	2.47	0.49
2:1E:237:ALA:O	2:1E:239:VAL:HG23	2.12	0.49
26:1H:1111:A:N3	26:1H:1112:G:H1'	2.27	0.49
9:82:44:VAL:O	9:82:51:ARG:NH2	2.45	0.49
26:1H:2650:U:H2'	26:1H:2651:C:C6	2.48	0.49
26:1H:2294:C:H2'	26:1H:2295:C:H6	1.78	0.49
1:1G:830:G:H2'	1:1G:831:U:O4'	2.12	0.49
27:16:90:C:P	37:88:16:ARG:HH21	2.35	0.49
1:1G:67:C:H2'	1:1G:68:G:C8	2.46	0.49
26:1H:1669:A:H5''	26:1H:1670:C:OP2	2.13	0.49
26:1H:475:U:C4	26:1H:481:G:O6	2.65	0.49
26:1H:2725:A:C4	26:1H:2727:G:C8	3.01	0.49
1:13:403:C:OP1	4:3E:137:SER:OG	2.30	0.49
1:1G:648:A:H2'	1:1G:649:G:H8	1.77	0.49
26:14:39:C:H2'	26:14:40:C:C6	2.48	0.49
30:39:46:ARG:HG2	30:39:46:ARG:NH1	2.23	0.49
26:1H:2378:A:H4'	39:A8:23:ARG:NH1	2.28	0.49
1:13:1148:U:H2'	1:13:1149:C:O4'	2.13	0.49
34:15:39:ARG:HD3	34:15:48:MET:HE2	1.95	0.49
1:1G:1273:G:H3'	1:1G:1274:G:C8	2.48	0.49
26:1H:2579:C:H2'	26:1H:2580:U:O4'	2.13	0.49
2:1E:162:ILE:HD11	2:1E:184:VAL:HG22	1.94	0.49
33:61:69:LYS:O	33:61:73:GLU:HB2	2.13	0.49
28:11:17:THR:HG22	28:11:204:ILE:HA	1.95	0.49
52:J5:49:CYS:SG	52:J5:50:GLY:N	2.86	0.49
26:1H:993:G:C4	26:1H:994:C:C5	3.01	0.49
1:1G:1517:G:C6	1:1G:1518:A:C5	3.00	0.49
26:14:27:G:O2'	26:14:28:A:OP2	2.28	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:1442:G:H2'	26:14:1443:G:H8	1.75	0.49
3:2E:167:TRP:CD1	3:2E:168:ALA:N	2.80	0.49
27:16:29:A:OP2	39:A8:31:SER:HB2	2.13	0.49
38:98:60:LEU:O	38:98:64:ARG:HG3	2.13	0.49
1:13:292:G:N7	1:13:293:G:H1'	2.27	0.49
26:1H:1808:U:H2'	26:1H:1809:A:O4'	2.12	0.49
26:14:2036:C:N4	58:14:3845:HOH:O	2.44	0.49
1:13:486:U:H2'	1:13:487:A:H8	1.77	0.49
20:BI:63:ILE:HG22	20:BI:77:ALA:HB1	1.94	0.49
1:1G:1051:C:H2'	1:1G:1052:U:C6	2.48	0.49
18:9A:45:SER:OG	18:9A:46:GLU:N	2.46	0.49
1:13:191(F):U:H2'	1:13:191:G:C8	2.47	0.49
26:14:566:U:O3'	58:14:3841:HOH:O	2.19	0.49
36:78:62:LEU:O	55:Q8:13:ARG:HD3	2.13	0.49
47:E5:48:GLY:HA3	47:E5:80:HIS:ND1	2.28	0.49
26:14:2012:G:O3'	43:A5:96:ILE:HG12	2.13	0.49
1:1G:854:G:C2	1:1G:855:G:C8	3.01	0.49
43:A5:13:SER:O	43:A5:16:LYS:HB2	2.13	0.49
16:7I:38:TYR:CE1	16:7I:50:LYS:HB2	2.47	0.49
26:14:1000:A:C6	26:14:1001:A:N1	2.81	0.49
26:1H:242:G:H5'	55:Q8:61:LEU:HD13	1.93	0.49
28:11:136:ILE:HG22	28:11:137:PRO:HD2	1.95	0.49
26:14:2207:C:O2	28:19:151:LYS:NZ	2.38	0.49
26:14:1364:G:N7	48:F5:2:SER:HB2	2.28	0.49
32:51:8:PRO:HG2	32:51:69:ARG:HH21	1.75	0.49
1:13:1015:A:H2'	1:13:1016:A:C8	2.48	0.49
46:D5:91:LEU:HB3	46:D5:130:PRO:HG3	1.94	0.49
8:72:12:ARG:CZ	8:72:27:PRO:HD3	2.42	0.49
3:22:156:ARG:HB3	3:22:160:ALA:O	2.13	0.49
1:1G:552:U:O2'	12:3A:86:ARG:O	2.26	0.49
33:69:88:ILE:HG22	33:69:90:GLY:H	1.78	0.49
28:19:85:ASP:OD2	28:19:88:ARG:HD2	2.13	0.49
2:1E:226:ARG:HG3	2:1E:227:GLY:H	1.77	0.49
11:2A:13:GLN:HA	11:2A:75:TYR:O	2.12	0.49
26:1H:90:U:H4'	26:1H:91:A:H5'	1.95	0.49
1:1G:607:A:H2'	1:1G:608:A:O4'	2.13	0.49
31:49:120:LEU:HG	31:49:179:PRO:O	2.13	0.49
26:1H:2280:G:C2'	26:1H:2281:C:H5'	2.42	0.49
1:1G:973:G:O6	1:1G:974:A:N6	2.46	0.48
19:AI:41:VAL:H	19:AI:44:MET:HB2	1.78	0.48
26:14:2331:G:H4'	47:E5:43:THR:N	2.28	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
30:31:78:ILE:HA	30:31:83:PHE:CD2	2.49	0.48
28:11:145:VAL:HB	28:11:155:LEU:HB2	1.95	0.48
29:29:119:ARG:HA	29:29:160:TYR:CD2	2.48	0.48
1:1G:1262:C:H2'	1:1G:1263:C:C6	2.48	0.48
28:11:131:LEU:HB2	28:11:136:ILE:HD11	1.95	0.48
7:6E:91:VAL:HG12	7:6E:95:ARG:HB3	1.95	0.48
1:1G:861:G:N2	1:1G:872:A:H2	2.11	0.48
1:1G:1095:U:H5''	1:1G:1109:C:O2	2.12	0.48
1:13:428:G:C8	1:13:430:A:C5	3.01	0.48
2:1E:17:PHE:H	2:1E:17:PHE:HD1	1.61	0.48
26:1H:2315:G:H5''	26:1H:2316:C:OP2	2.13	0.48
1:13:1060:C:C5	3:2E:2:GLY:HA2	2.47	0.48
30:31:130:ALA:H	30:31:132:VAL:HG13	1.78	0.48
26:14:1794:U:H2'	26:14:1795:C:H6	1.77	0.48
1:13:711:G:O2'	1:13:712:A:H5'	2.13	0.48
1:13:313:A:H2'	1:13:314:C:H6	1.78	0.48
26:1H:1591:G:H2'	26:1H:1592:C:H6	1.78	0.48
29:21:73:GLU:HG3	29:21:74:PRO:HD2	1.95	0.48
7:6E:45:ASP:O	7:6E:49:ILE:HG12	2.13	0.48
26:14:1316:U:H2'	26:14:1317:A:C8	2.48	0.48
7:62:62:PHE:HD1	7:62:124:LEU:HD11	1.77	0.48
26:1H:818:G:H4'	26:1H:838:C:O3'	2.13	0.48
26:14:1337:G:H2'	26:14:1338:G:H8	1.78	0.48
29:29:113:PHE:HA	29:29:159:HIS:HD2	1.78	0.48
51:I5:43:TYR:O	51:I5:43:TYR:CG	2.66	0.48
38:55:78:LYS:O	38:55:83:ILE:HG13	2.13	0.48
26:1H:2850:A:C2	26:1H:2851:A:C4	3.01	0.48
5:4E:29:GLY:HA2	5:4E:46:GLY:O	2.12	0.48
1:1G:60:A:N6	1:1G:110:C:N3	2.59	0.48
26:1H:2791:C:H42	26:1H:2805:G:H1	1.61	0.48
1:1G:868:C:H2'	1:1G:869:G:O4'	2.12	0.48
20:BA:75:ASN:N	20:BA:75:ASN:OD1	2.45	0.48
1:13:12:U:O2'	1:13:526:C:H4'	2.13	0.48
12:3I:39:VAL:HG12	12:3I:41:ARG:HG2	1.95	0.48
1:1G:1509:C:H2'	1:1G:1510:U:O4'	2.12	0.48
1:13:545:C:O2'	1:13:549:C:OP1	2.30	0.48
42:D8:3:ALA:HB1	42:D8:38:LEU:HD11	1.94	0.48
1:1G:1499:A:H1'	1:1G:1520:G:H5'	1.94	0.48
1:1G:19:C:H5''	5:42:86:ALA:HB1	1.95	0.48
10:1I:90:LEU:N	10:1I:91:PRO:HD3	2.28	0.48
26:1H:1576:U:N3	26:1H:1577:C:C5	2.81	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1G:1256:A:H2'	1:1G:1278:U:O2	2.13	0.48
47:E5:18:ALA:CB	47:E5:20:ARG:HE	2.26	0.48
26:14:273(D):C:N4	26:14:363(B):G:H1	2.11	0.48
38:98:78:LYS:O	38:98:83:ILE:HG13	2.13	0.48
41:85:90:VAL:HA	42:95:39:LEU:HD23	1.95	0.48
4:32:59:ARG:HA	4:32:62:GLN:HB2	1.95	0.48
31:41:173:LEU:HD22	31:41:178:PHE:CE2	2.48	0.48
26:14:1210:A:H5'	26:14:1212:G:O4'	2.13	0.48
23:2K:57:C:H5''	23:2K:58:A:OP2	2.13	0.48
1:1G:1131:G:C8	1:1G:1132:C:H5	2.31	0.48
46:D5:138:GLU:O	46:D5:156:LYS:HG3	2.13	0.48
35:68:73:ASP:OD1	35:68:75:SER:N	2.46	0.48
1:13:31:G:HO2'	1:13:48:C:N4	2.10	0.48
26:1H:363(A):A:H2'	26:1H:363(B):G:H8	1.77	0.48
9:82:21:PRO:HA	9:82:59:PHE:HD1	1.78	0.48
1:13:392:G:H5'	16:7I:12:LYS:HE3	1.95	0.48
8:7E:104:ARG:O	8:7E:107:LEU:HB2	2.12	0.48
30:39:111:ALA:HB2	30:39:206:ILE:HG21	1.94	0.48
26:14:1430:C:H2'	26:14:1431:U:C6	2.49	0.48
53:O8:9:LEU:HB3	53:O8:26:ASN:O	2.13	0.48
33:69:2:LYS:HA	33:69:20:ASP:HA	1.95	0.48
26:14:1859:A:N6	26:14:1883:G:O2'	2.46	0.48
50:H5:12:PRO:HB2	50:H5:20:LYS:HG3	1.95	0.48
26:1H:845:G:H8	26:1H:845:G:OP2	1.96	0.48
33:61:77:LEU:HD23	33:61:101:LEU:HB2	1.96	0.48
51:M8:15:ILE:HD11	51:M8:32:TYR:CD1	2.48	0.48
41:C8:43:GLY:HA3	42:D8:73:SER:OG	2.13	0.48
51:I5:27:THR:HG22	51:I5:28:LYS:HG3	1.94	0.48
30:39:20:LEU:HG	30:39:199:TRP:HH2	1.77	0.48
29:29:63:LEU:HD23	29:29:63:LEU:H	1.77	0.48
9:8E:112:LYS:CA	9:8E:119:ALA:HB2	2.41	0.48
1:13:923:A:O2'	1:13:1399:C:OP2	2.23	0.48
49:G5:32:LEU:HB2	49:G5:53:LEU:HD13	1.93	0.48
26:1H:1728:G:C6	26:1H:1730:U:OP2	2.67	0.48
26:1H:1385:G:O2'	26:1H:1396:U:C6	2.62	0.48
24:3K:56:C:N4	26:1H:2112:G:H1	2.11	0.48
30:39:117:ARG:HD3	30:39:120:GLU:OE1	2.13	0.48
26:14:2748:A:H2'	26:14:2749:A:H8	1.78	0.48
26:1H:356:G:H2'	26:1H:357:A:C8	2.47	0.48
26:14:35:G:C4	26:14:454:A:C2	3.02	0.48
29:21:13:ARG:HD2	40:B8:58:ASN:HB3	1.94	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
11:2I:32:ILE:HD12	11:2I:72:ALA:HB2	1.95	0.48
2:12:21:ARG:HA	2:12:39:ILE:HA	1.94	0.48
26:14:200:U:O2	26:14:386:G:N2	2.47	0.48
24:3L:30:G:N2	24:3L:41:C:N3	2.62	0.48
55:Q8:31:HIS:CD2	55:Q8:34:TRP:CH2	3.02	0.48
26:14:601:C:O2	26:14:605:C:H4'	2.13	0.48
1:13:55:A:H2	33:69:82:ARG:HG2	1.79	0.48
29:29:120:TRP:HE1	29:29:156:MET:H	1.61	0.48
26:1H:197:A:N6	26:1H:2430:A:H2'	2.27	0.48
41:C8:106:PHE:O	41:C8:109:LEU:HB2	2.14	0.48
26:1H:1204:A:C2	26:1H:1241:A:N1	2.82	0.48
1:1G:1469:G:H2'	1:1G:1470:G:C8	2.48	0.48
26:14:753:C:O2'	26:14:754:C:H5'	2.14	0.48
26:1H:1252:G:H5''	58:1H:3993:HOH:O	2.12	0.48
46:D5:9:TYR:HD1	46:D5:37:VAL:HG12	1.78	0.48
32:51:20:ALA:HB1	32:51:21:PRO:HD2	1.95	0.48
26:14:1798:U:H5'	28:19:259:THR:OG1	2.14	0.48
41:85:106:PHE:O	41:85:109:LEU:HB2	2.13	0.48
37:88:101:ARG:HG3	37:88:102:VAL:N	2.29	0.48
55:M5:33:ASN:O	55:M5:36:LYS:NZ	2.36	0.48
34:15:95:PRO:O	34:15:98:VAL:HG22	2.13	0.48
26:14:807:U:H2'	26:14:808:G:H8	1.77	0.48
1:13:186(A):C:O2	20:BI:104:LEU:HD12	2.14	0.48
26:1H:2093:G:OP2	33:61:22:LYS:HD2	2.13	0.48
42:95:69:LYS:HE3	42:95:85:LYS:HD2	1.96	0.48
26:14:1486:A:H2'	26:14:1487:G:C8	2.48	0.48
26:1H:2061:G:P	58:1H:3632:HOH:O	2.70	0.48
28:11:85:ASP:OD2	28:11:88:ARG:NH1	2.41	0.48
10:1I:49:VAL:HG23	14:5I:41:ARG:HD2	1.94	0.48
26:1H:1142:U:H5'	26:1H:1142(A):A:C8	2.48	0.48
1:1G:430:A:OP2	4:32:8:VAL:HG23	2.14	0.48
26:14:933:A:C5	26:14:934:G:C8	3.01	0.48
13:4I:94:ARG:HH22	26:1H:887:A:C5'	2.26	0.48
15:6I:70:LEU:HD11	15:6I:77:ARG:HG3	1.94	0.48
13:4A:3:ARG:NH1	13:4A:7:VAL:HG12	2.28	0.48
39:A8:83:LYS:HG3	39:A8:110:LEU:HD12	1.95	0.48
23:2L:41:C:C2	23:2L:42:C:C5	3.02	0.48
1:13:648:A:C6	1:13:649:G:C6	3.01	0.48
38:98:9:LYS:HA	38:98:17:ARG:HE	1.77	0.48
3:22:73:PRO:O	3:22:77:ILE:HG13	2.13	0.48
26:1H:2176:A:H2'	26:1H:2177:C:C6	2.48	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:508:G:N2	43:A5:80:PRO:HG2	2.27	0.48
38:55:28:LEU:HD12	38:55:48:VAL:HG21	1.95	0.48
45:G8:27:VAL:HA	45:G8:39:VAL:HG22	1.95	0.48
43:E8:110:LYS:HG3	43:E8:111:HIS:H	1.79	0.48
26:1H:301:G:C4	26:1H:302:C:C5	3.02	0.48
26:14:1849:G:H2'	26:14:1850:G:C8	2.48	0.48
4:32:191:ARG:NH2	4:32:194:LEU:O	2.46	0.48
1:13:1266:G:N2	1:13:1270:C:C2	2.81	0.48
26:1H:190:A:OP2	48:J8:39:LYS:HE3	2.13	0.48
37:45:2:LEU:O	37:45:70:PRO:HG2	2.12	0.48
26:1H:140:A:H8	26:1H:1408:C:O2'	1.94	0.48
26:1H:2376:A:C2	39:A8:112:PHE:HB3	2.48	0.48
46:H8:15:PRO:O	46:H8:19:ARG:HB2	2.13	0.48
46:H8:75:ASN:O	46:H8:84:GLU:N	2.36	0.48
26:1H:1790:C:H2'	26:1H:1791:A:C5	2.48	0.48
26:14:1071:G:N2	26:14:1087:G:H22	2.12	0.48
1:1G:1187:G:H4'	9:82:111:ARG:NH1	2.28	0.48
40:75:25:GLY:H	40:75:49:VAL:HG23	1.78	0.48
26:1H:2080:G:C8	26:1H:2080:G:H5''	2.47	0.48
26:14:2297:C:H2'	26:14:2298:A:C8	2.46	0.48
26:14:1278:A:C5'	38:55:36:THR:HG22	2.42	0.48
26:1H:298:G:N7	58:1H:4010:HOH:O	2.35	0.48
41:85:88:ILE:HA	42:95:49:THR:O	2.13	0.48
1:1G:79:G:N2	1:1G:90:C:N3	2.58	0.48
27:1J:12:C:O2'	47:E5:74:ARG:HG2	2.13	0.48
1:1G:1112:C:C4	3:22:178:LEU:HD23	2.48	0.48
26:1H:2619:C:O2'	26:1H:2620:C:H5'	2.12	0.48
26:14:1889:A:O2'	26:14:2087:G:H5'	2.13	0.48
26:14:2746:U:OP1	32:59:85:LYS:NZ	2.46	0.48
30:31:149:ASP:OD1	30:31:149:ASP:N	2.44	0.48
26:14:1530:G:H2'	26:14:1531:C:O4'	2.13	0.48
26:14:533:G:H2'	26:14:534:U:O4'	2.14	0.48
27:1J:63:G:C2	27:1J:64:C:C2	3.02	0.48
55:Q8:46:ARG:HH21	55:Q8:48:PHE:HB2	1.78	0.48
26:1H:71:A:H2	44:F8:31:HIS:NE2	2.04	0.48
26:14:2749:A:N1	26:14:2750:A:N6	2.61	0.48
26:1H:1465:G:C4	26:1H:1466:G:C8	3.02	0.48
11:2A:86:GLY:H	11:2A:112:THR:HG23	1.79	0.48
8:7E:86:ILE:HG22	8:7E:87:SER:N	2.28	0.48
8:7E:87:SER:CB	8:7E:93:VAL:H	2.26	0.48
1:1G:1060:C:O2'	10:1A:56:HIS:ND1	2.45	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
30:39:7:TYR:HD1	30:39:18:ARG:H	1.61	0.48
9:82:10:ARG:HA	9:82:104:ARG:HH21	1.79	0.48
7:6E:92:SER:O	7:6E:96:GLN:HG3	2.12	0.48
1:13:160:A:H61	1:13:347:G:H1'	1.79	0.48
1:13:590:C:H42	1:13:649:G:H1	1.60	0.48
26:14:747:U:O2	26:14:2014:A:H1'	2.14	0.48
29:21:37:ARG:HA	29:21:42:ASP:OD2	2.14	0.48
12:3A:36:VAL:HG13	12:3A:81:SER:H	1.79	0.48
18:9A:31:LEU:H	18:9A:31:LEU:HD23	1.79	0.48
26:1H:2331:G:H4'	47:18:42:GLY:HA3	1.95	0.48
26:1H:195:A:H4'	26:1H:251:A:O2'	2.13	0.48
23:2L:32:G:C5	23:2L:33:OMC:C5	3.01	0.48
1:13:284:G:H2'	1:13:285:G:H8	1.79	0.48
34:58:6:PRO:HG3	34:58:41:ASP:HB2	1.94	0.48
6:52:15:ASP:OD1	6:52:17:SER:N	2.46	0.48
34:58:99:LEU:HD23	34:58:99:LEU:HA	1.74	0.48
48:J8:64:ALA:HA	48:J8:67:ILE:HG13	1.95	0.48
1:13:1053:G:O5'	1:13:1054:C:H3'	2.14	0.48
31:49:76:SER:OG	31:49:84:LYS:N	2.47	0.48
35:68:118:ALA:HA	35:68:119:PRO:HD2	1.75	0.48
26:1H:2615:U:H2'	26:1H:2616:C:H6	1.79	0.48
1:1G:321:A:N7	1:1G:328:C:C6	2.81	0.48
39:65:67:ARG:HB2	39:65:67:ARG:CZ	2.43	0.48
1:1G:397:A:N3	1:1G:397:A:H3'	2.28	0.48
26:14:1485:G:H2'	26:14:1486:A:H8	1.77	0.48
26:14:38:A:H2'	26:14:39:C:C6	2.49	0.48
50:L8:5:LYS:HD2	50:L8:34:GLU:OE1	2.14	0.48
1:13:1026:G:H1	1:13:1035:A:H61	1.62	0.48
26:1H:1021:A:C8	26:1H:1022:G:H5''	2.48	0.48
26:1H:1676:A:C2	26:1H:1993:U:H5'	2.48	0.48
26:1H:960:A:H61	37:88:82:ARG:HH21	1.62	0.48
29:29:51:PHE:CE2	29:29:52:LEU:HG	2.48	0.48
36:35:80:TYR:CD2	36:35:111:ARG:HB3	2.49	0.48
26:1H:993:G:H5''	41:C8:50:ARG:HH21	1.78	0.48
26:1H:99:U:C6	26:1H:102:G:C2	3.01	0.48
12:3A:47:LYS:O	12:3A:49:ASN:N	2.46	0.48
28:19:83:GLU:OE1	28:19:104:TYR:OH	2.23	0.48
33:61:95:LYS:HE3	33:61:99:GLU:CD	2.34	0.48
18:9A:36:ASN:ND2	18:9A:39:VAL:HB	2.29	0.48
46:D5:28:MET:O	46:D5:35:ARG:N	2.34	0.48
26:1H:1170:G:N2	26:1H:1180:C:C2	2.81	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:1653:G:C6	38:55:9:LYS:HB2	2.49	0.48
26:1H:1543:A:H3'	26:1H:1543:A:OP2	2.13	0.48
29:21:92:THR:HB	29:21:94:GLU:HG2	1.95	0.48
37:88:39:PRO:HA	37:88:97:VAL:O	2.13	0.48
26:1H:487:C:H2'	26:1H:488:G:H5'	1.94	0.48
26:14:1003:G:N2	26:14:1153:C:C2	2.81	0.48
44:B5:57:LEU:N	44:B5:57:LEU:HD23	2.29	0.48
39:65:28:VAL:HG11	39:65:98:VAL:HG12	1.94	0.48
26:14:1354:A:OP2	58:14:3772:HOH:O	2.20	0.48
36:78:85:LEU:HD13	36:78:120:ALA:HB2	1.95	0.48
1:13:465:A:N7	1:13:467:G:C5	2.82	0.48
1:1G:1126:U:C4	1:1G:1281:U:C6	3.02	0.48
26:1H:900:A:H5'	26:1H:901:A:OP2	2.14	0.48
26:1H:783:A:C8	26:1H:783:A:H3'	2.48	0.48
26:14:2872:G:C4	26:14:2873:A:C2	3.01	0.48
26:1H:1533:C:H2'	26:1H:1534:G:C2	2.49	0.48
1:1G:942:G:N2	9:82:124:GLN:OE1	2.35	0.48
16:7I:28:ARG:HG3	16:7I:29:ASP:N	2.28	0.48
26:14:972:G:OP2	26:14:974:G:H5''	2.13	0.48
2:12:19:HIS:HE1	2:12:206:ASP:HB2	1.79	0.48
26:1H:107:C:H2'	26:1H:108:U:C6	2.49	0.48
42:D8:25:LEU:H	42:D8:92:THR:HG23	1.79	0.48
1:1G:1338:G:C6	1:1G:1339:A:N1	2.82	0.48
20:BA:67:ALA:HA	20:BA:73:HIS:HA	1.94	0.48
46:H8:53:ILE:HG22	46:H8:71:VAL:HG22	1.95	0.48
1:13:486:U:H2'	1:13:487:A:C8	2.48	0.48
4:3E:7:PRO:HB2	4:3E:10:ARG:HG2	1.95	0.48
1:1G:667:G:N2	1:1G:739:C:O2	2.42	0.48
47:I8:50:ASN:ND2	47:I8:83:PRO:HD3	2.29	0.48
33:61:93:THR:N	33:61:96:ASP:OD2	2.38	0.48
26:1H:207:A:H2'	26:1H:208:C:O4'	2.14	0.48
26:1H:1491:G:H5'	28:11:99:ASP:OD1	2.14	0.48
26:14:1418:G:H8	26:14:1418:G:O5'	1.96	0.48
28:19:67:PHE:HB3	28:19:153:ALA:H	1.78	0.48
5:4E:80:ILE:HG13	5:4E:81:GLU:N	2.28	0.48
40:B8:50:ILE:HG22	40:B8:62:THR:OG1	2.14	0.48
1:1G:1149:C:OP2	9:82:9:ARG:NH1	2.47	0.48
26:1H:1264:G:OP1	52:N8:19:ARG:NH2	2.27	0.48
5:4E:63:ARG:HA	5:4E:66:MET:HE2	1.96	0.48
50:L8:54:VAL:HG13	50:L8:54:VAL:O	2.14	0.48
1:1G:1220:G:O3'	19:AA:36:ARG:HD3	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:2439:A:C5'	26:14:2439:A:H8	2.27	0.48
23:2K:54:G:H2'	23:2K:55:5MU:C6	2.43	0.48
26:1H:2496:C:OP1	37:88:82:ARG:HD3	2.14	0.48
1:13:101:A:OP2	1:13:101:A:H8	1.97	0.48
26:14:2295:C:OP2	39:65:10:ARG:HD3	2.13	0.48
37:45:57:HIS:CD2	37:45:117:ALA:HB2	2.49	0.48
26:14:1427:A:H4'	26:14:1428:C:O5'	2.14	0.48
17:8I:76:LEU:HD21	17:8I:79:SER:HB3	1.95	0.48
26:1H:910:A:H62	37:88:12:GLN:HA	1.79	0.48
1:1G:1218:C:OP2	14:5A:9:LYS:NZ	2.37	0.48
26:1H:1794:U:H2'	26:1H:1795:C:H6	1.78	0.48
30:39:78:ILE:HA	30:39:83:PHE:CD2	2.48	0.48
1:13:983:A:H5'	14:5I:3:ARG:HH12	1.79	0.48
26:1H:1221:C:H2'	26:1H:1222:C:C6	2.48	0.48
47:18:38:VAL:CG1	47:18:40:GLN:HG2	2.44	0.48
3:2E:62:ASP:N	3:2E:62:ASP:OD1	2.47	0.48
17:8I:74:LEU:HA	17:8I:74:LEU:HD22	1.74	0.48
37:45:36:ALA:HB2	37:45:103:MET:SD	2.54	0.48
1:1G:17:U:H2'	1:1G:18:C:C6	2.49	0.48
45:G8:5:MET:HG3	45:G8:6:HIS:H	1.78	0.48
1:1G:1468:A:H5''	1:1G:1469:G:OP2	2.12	0.48
34:15:135:PRO:HB2	34:15:137:LYS:HZ1	1.78	0.48
13:4I:50:GLU:HA	13:4I:53:VAL:HB	1.95	0.48
13:4I:50:GLU:O	13:4I:54:VAL:HG23	2.13	0.48
37:88:75:THR:HA	37:88:87:LYS:O	2.13	0.48
16:7I:19:ILE:HB	16:7I:36:ILE:O	2.14	0.48
1:1G:195:A:N7	1:1G:196:A:C6	2.82	0.48
32:51:101:ARG:NH1	32:51:122:THR:OG1	2.31	0.48
44:F8:24:GLY:O	44:F8:83:VAL:HG22	2.14	0.48
16:7A:28:ARG:NE	16:7A:29:ASP:OD1	2.37	0.48
40:75:118:ARG:NH1	40:75:121:ILE:HG21	2.29	0.48
26:1H:2239:G:H5'	28:11:251:GLY:HA3	1.96	0.48
26:1H:2050:C:H2'	26:1H:2051:A:O4'	2.14	0.48
26:14:491:G:O6	43:A5:49:LYS:HD3	2.14	0.48
1:13:1366:C:H2'	1:13:1367:C:C6	2.49	0.48
26:1H:1730:U:H5''	26:1H:1731:G:OP2	2.14	0.48
1:1G:1298:C:C5	7:62:114:ARG:HD2	2.48	0.48
26:1H:654(M):C:H5'	26:1H:654(N):G:C5	2.49	0.48
1:13:963:G:H1	1:13:972:C:N4	2.10	0.48
1:13:1073:U:O2'	2:1E:104:ASN:ND2	2.37	0.48
26:1H:1550:C:H2'	26:1H:1551:C:C6	2.49	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
38:55:38:VAL:HG22	38:55:112:ALA:HB2	1.95	0.48
26:1H:2318:G:N2	39:A8:2:ALA:HA	2.29	0.48
1:1G:1238:A:N3	1:1G:1241:G:O2'	2.28	0.48
26:1H:2287:A:N1	26:1H:2346:A:H2	2.12	0.48
26:1H:2306:C:H3'	26:1H:2307:G:C5'	2.43	0.48
26:1H:583:G:H5''	41:C8:10:ARG:HH12	1.78	0.48
26:1H:1705:G:H2'	26:1H:1706:U:H5'	1.96	0.48
33:61:69:LYS:HG3	33:61:136:VAL:HB	1.96	0.48
8:7E:9:MET:HG3	8:7E:26:VAL:HG21	1.96	0.48
49:K8:31:GLU:HB2	49:K8:53:LEU:HD11	1.95	0.48
26:1H:1001:A:C8	26:1H:1002:G:C8	3.02	0.48
24:3L:23:A:H2'	24:3L:24:G:C8	2.49	0.48
1:1G:789:U:O2'	1:1G:791:G:O6	2.32	0.48
1:1G:1386:G:H2'	1:1G:1387:G:H8	1.79	0.48
26:14:2763:G:OP2	58:14:4000:HOH:O	2.20	0.48
43:A5:70:TYR:O	43:A5:107:LEU:HD12	2.14	0.48
26:14:2593:U:H2'	26:14:2594:C:C6	2.49	0.48
26:1H:2756:U:H4'	26:1H:2757:A:OP1	2.13	0.48
1:1G:1291:G:H4'	9:82:39:GLY:HA3	1.95	0.48
26:14:673:C:H4'	30:39:82:ILE:CG1	2.44	0.48
31:49:72:ARG:HD2	31:49:85:GLY:O	2.13	0.48
46:D5:127:LYS:O	46:D5:162:GLU:HB2	2.14	0.48
26:1H:1831:G:H2'	26:1H:1832:C:C6	2.49	0.48
26:14:2290:G:O2'	26:14:2381:C:H1'	2.13	0.48
1:13:1347:G:H5''	9:8E:107:ARG:HG2	1.95	0.47
26:14:2420:C:N4	55:M5:31:HIS:HB3	2.29	0.47
21:1B:9:ARG:HG3	21:1B:10:ARG:N	2.29	0.47
29:29:13:ARG:NH2	40:75:77:PRO:HB3	2.29	0.47
1:1G:660:G:N2	1:1G:745:C:N3	2.53	0.47
26:14:1379:A:H4'	26:14:1380:G:OP2	2.14	0.47
4:3E:18:LYS:HD3	4:3E:31:CYS:SG	2.54	0.47
1:1G:1255:G:O3'	1:1G:1258:G:H1'	2.14	0.47
1:1G:1240:U:H2'	7:62:32:ARG:NH2	2.29	0.47
42:D8:43:GLU:OE2	42:D8:44:LYS:HG2	2.14	0.47
26:14:869:G:C2	26:14:870:A:C8	3.02	0.47
26:14:5:A:H2'	26:14:6:A:H8	1.78	0.47
26:1H:2400:G:H2'	26:1H:2401:U:C6	2.48	0.47
26:1H:1337:G:H2'	26:1H:1338:G:C8	2.48	0.47
37:45:17:LEU:HD21	37:45:41:TRP:NE1	2.28	0.47
5:4E:150:ARG:CZ	5:4E:150:ARG:HB3	2.44	0.47
10:1A:78:ASN:OD1	10:1A:80:LYS:HB3	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1G:1453:G:H8	20:BA:39:LYS:HD2	1.79	0.47
5:42:84:PHE:O	5:42:86:ALA:N	2.37	0.47
1:13:1113:C:H2'	1:13:1114:C:H6	1.79	0.47
1:13:1236:A:O2'	1:13:1304:G:H4'	2.14	0.47
54:P8:10:ARG:HG2	54:P8:10:ARG:O	2.14	0.47
14:5A:4:LYS:O	14:5A:7:ILE:HG12	2.13	0.47
44:B5:84:ALA:O	44:B5:87:GLN:HG3	2.13	0.47
28:19:267:SER:O	28:19:268:ARG:HG2	2.14	0.47
1:1G:762:C:H2'	1:1G:763:G:H8	1.78	0.47
49:K8:8:LYS:HZ2	49:K8:12:GLU:HG2	1.79	0.47
1:1G:604:G:H2'	1:1G:605:U:O4'	2.14	0.47
26:14:2353:G:H2'	26:14:2354:G:O4'	2.13	0.47
1:13:985:C:H2'	1:13:986:A:C8	2.49	0.47
1:13:34:C:H2'	1:13:35:G:C8	2.49	0.47
1:13:825:G:O4'	8:7E:2:LEU:HD21	2.14	0.47
3:22:38:ARG:O	3:22:42:LEU:N	2.31	0.47
1:1G:801:U:H2'	1:1G:802:A:C8	2.49	0.47
38:55:104:ARG:HD2	38:55:107:ASP:OD2	2.13	0.47
2:12:5:ILE:HG12	2:12:6:THR:HG22	1.95	0.47
1:1G:1154:G:N3	1:1G:1155:G:C8	2.83	0.47
26:1H:1408:C:C2	26:1H:1595:G:N2	2.82	0.47
26:1H:654(D):G:H22	26:1H:654(Q):C:N4	2.12	0.47
26:1H:1442:G:C2	26:1H:1550:C:O2	2.67	0.47
38:55:38:VAL:HB	38:55:39:PRO:HD3	1.97	0.47
24:1L:18:G:O2'	24:1L:60:U:O4	2.26	0.47
29:29:89:ASP:CG	29:29:90:THR:H	2.17	0.47
19:AA:15:LEU:HD12	19:AA:33:THR:HB	1.95	0.47
34:15:15:LEU:HD13	34:15:16:ILE:N	2.28	0.47
1:1G:1240:U:OP2	7:62:116:ALA:N	2.43	0.47
7:62:116:ALA:HA	7:62:119:ARG:HE	1.79	0.47
41:85:91:ASP:CG	41:85:96:ALA:HB2	2.35	0.47
4:32:108:LEU:CD1	4:32:174:LEU:HB3	2.44	0.47
26:14:2065:C:H2'	26:14:2066:C:H6	1.79	0.47
37:88:21:THR:HG22	37:88:99:PRO:O	2.13	0.47
29:21:70:ALA:O	29:21:73:GLU:N	2.47	0.47
26:14:1316:U:H2'	26:14:1317:A:H8	1.79	0.47
26:1H:2562:U:H1'	35:68:23:ARG:HD3	1.95	0.47
26:1H:2022:U:O2'	26:1H:2617:C:H5'	2.14	0.47
1:13:186(E):C:N3	1:13:191(C):G:C2	2.82	0.47
26:1H:2785:C:OP1	29:21:41:LYS:HE3	2.14	0.47
26:14:1259:G:H2'	26:14:1260:G:H8	1.79	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
45:C5:83:THR:HG22	45:C5:84:ARG:H	1.79	0.47
26:1H:2238:G:H4'	26:1H:2239:G:OP1	2.14	0.47
1:1G:578:C:H42	1:1G:763:G:H1	1.60	0.47
1:13:838:G:H1	1:13:848:C:N4	2.12	0.47
1:1G:1459:C:OP1	20:BA:31:SER:OG	2.30	0.47
26:14:2861:G:O2'	26:14:2862:G:H5'	2.14	0.47
12:3A:70:ILE:HD13	12:3A:77:LEU:HD12	1.95	0.47
2:12:215:LEU:HA	2:12:218:ALA:HB3	1.97	0.47
26:14:2619:C:OP1	29:29:152:LYS:HE3	2.14	0.47
32:59:99:VAL:HG13	32:59:100:GLY:H	1.78	0.47
10:1A:36:GLY:O	10:1A:38:ILE:HG13	2.13	0.47
1:1G:308:C:H2'	1:1G:309:G:C8	2.49	0.47
10:1A:9:ARG:NH2	10:1A:95:GLU:OE1	2.46	0.47
1:1G:790:A:O5'	1:1G:790:A:H8	1.97	0.47
1:1G:1280:A:H5'	1:1G:1281:U:OP2	2.14	0.47
24:1L:74:C:O2'	24:1L:75:C:OP2	2.28	0.47
26:1H:446:G:OP2	58:1H:3691:HOH:O	2.19	0.47
1:1G:616:G:C2	1:1G:617:G:C8	3.02	0.47
1:1G:616:G:N3	1:1G:617:G:C8	2.82	0.47
1:1G:1037:C:H2'	1:1G:1038:C:C6	2.49	0.47
26:1H:1534:G:H22	26:1H:1538:G:N2	2.09	0.47
26:1H:1022:G:O6	34:58:66:LYS:NZ	2.47	0.47
26:14:2439:A:C5'	26:14:2439:A:C8	2.96	0.47
1:13:266:G:H5''	1:13:267:C:H5	1.78	0.47
37:88:81:VAL:O	37:88:82:ARG:HB2	2.14	0.47
28:11:231:HIS:CD2	28:11:249:PRO:HA	2.49	0.47
1:1G:1261:A:C2	1:1G:1262:C:H1'	2.50	0.47
13:4I:12:ASN:O	13:4I:45:VAL:HG12	2.13	0.47
48:F5:76:ARG:HG3	48:F5:94:LEU:HD13	1.95	0.47
3:22:44:GLU:HA	3:22:52:LEU:HD11	1.95	0.47
26:1H:1055:G:O2'	26:1H:1086:A:N6	2.47	0.47
26:14:861:A:C2	26:14:917:A:C4	3.03	0.47
24:3K:23:A:H2'	24:3K:24:G:C8	2.48	0.47
24:3K:9:A:N6	24:3K:23:A:H62	2.13	0.47
33:61:52:ARG:O	33:61:56:LYS:HB3	2.15	0.47
1:13:1326:C:H2'	1:13:1327:C:H6	1.80	0.47
3:22:112:SER:HB3	3:22:115:LEU:HD12	1.96	0.47
53:O8:47:THR:HG22	53:O8:48:VAL:HG23	1.96	0.47
6:5E:41:GLU:HB2	6:5E:62:TRP:HE3	1.80	0.47
1:1G:57:G:H2'	1:1G:58:C:C6	2.49	0.47
26:14:754:C:H2'	26:14:755:C:C6	2.49	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:1H:2383:G:O2'	26:1H:2384:G:H5'	2.13	0.47
26:1H:2591:C:P	28:11:239:ARG:HG3	2.54	0.47
1:1G:303:A:H2'	1:1G:304:U:O4'	2.14	0.47
45:C5:90:LEU:HA	45:C5:91:GLU:HA	1.74	0.47
46:H8:92:SER:O	46:H8:130:PRO:HG2	2.13	0.47
36:78:63:PRO:HG2	55:Q8:24:ALA:HB3	1.96	0.47
34:58:7:LYS:HB3	34:58:7:LYS:HE3	1.67	0.47
26:14:1358:G:N2	26:14:1372:U:C5	2.82	0.47
26:1H:330:A:O2'	26:1H:331:A:C8	2.58	0.47
26:1H:900:A:H3'	26:1H:901:A:C8	2.45	0.47
22:1K:76:A:H1'	26:1H:2583:G:H21	1.80	0.47
26:1H:1156:A:P	58:1H:3762:HOH:O	2.69	0.47
40:B8:88:ILE:HD13	40:B8:91:ARG:HH21	1.79	0.47
27:1J:3:C:H2'	27:1J:4:C:C6	2.49	0.47
26:1H:2376:A:H2'	26:1H:2377:A:O4'	2.14	0.47
16:7I:5:ARG:HH21	16:7I:22:THR:HG21	1.79	0.47
55:M5:34:TRP:HD1	55:M5:34:TRP:H	1.62	0.47
1:13:129(A):G:N2	1:13:188:U:O2'	2.48	0.47
26:1H:635:C:O2'	26:1H:639:U:OP1	2.32	0.47
1:13:692:U:O2'	1:13:694:A:N7	2.37	0.47
23:2K:62:C:O2'	23:2K:63:C:H5'	2.14	0.47
3:2E:91:LEU:O	3:2E:95:THR:OG1	2.29	0.47
23:2K:8:4SU:O2	23:2K:22:A:H2	1.97	0.47
26:14:1387:C:C2	26:14:1388:G:C8	3.01	0.47
26:1H:35:G:H2'	26:1H:36:G:O4'	2.13	0.47
7:62:65:ALA:HB2	7:62:128:ALA:HB2	1.96	0.47
18:9I:66:LEU:HD11	18:9I:70:ILE:HD11	1.95	0.47
26:1H:2074:U:P	58:1H:3682:HOH:O	2.72	0.47
26:14:192:C:OP1	58:14:3628:HOH:O	2.20	0.47
36:78:94:GLU:HG3	36:78:96:THR:HG23	1.97	0.47
45:C5:87:LYS:H	45:C5:94:LYS:HG2	1.79	0.47
39:A8:67:ARG:O	39:A8:71:ARG:HG3	2.15	0.47
37:88:16:ARG:HE	37:88:16:ARG:HB3	1.63	0.47
26:14:807:U:C2	26:14:808:G:C8	3.03	0.47
32:59:20:ALA:O	32:59:22:GLY:N	2.46	0.47
45:G8:20:TYR:CE1	45:G8:43:ASN:HA	2.49	0.47
30:39:118:ALA:O	30:39:121:GLY:N	2.46	0.47
10:1I:40:LEU:HB2	10:1I:69:ASN:HB2	1.96	0.47
26:14:506:G:O3'	26:14:507:A:H8	1.98	0.47
42:95:29:PRO:HA	42:95:61:VAL:HG11	1.95	0.47
32:51:129:THR:OG1	32:51:129:THR:O	2.25	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:1353:A:OP2	58:14:3768:HOH:O	2.20	0.47
31:41:82:LEU:HD21	31:41:88:ILE:HD11	1.97	0.47
26:1H:587:C:H4'	26:1H:588:U:H6	1.79	0.47
1:13:1346:A:OP1	9:8E:120:ARG:NH1	2.47	0.47
29:21:144:ARG:HG3	29:21:144:ARG:HH11	1.80	0.47
26:14:311:A:C6	26:14:328:U:C4	3.03	0.47
45:C5:40:GLU:N	45:C5:40:GLU:OE2	2.47	0.47
58:1H:4235:HOH:O	48:J8:65:SER:HA	2.13	0.47
26:14:898:C:H3'	26:14:899:A:H5''	1.95	0.47
8:7E:87:SER:HB3	8:7E:133:LEU:O	2.14	0.47
1:1G:1172:C:H2'	1:1G:1173:G:H8	1.79	0.47
1:13:232:G:C5	1:13:233:C:C5	3.02	0.47
4:3E:173:TRP:CG	4:3E:189:PRO:HG3	2.48	0.47
1:13:953:G:O5'	1:13:953:G:H8	1.97	0.47
26:14:1729:A:H2'	26:14:1731:G:H22	1.80	0.47
49:K8:27:GLU:O	49:K8:31:GLU:HG3	2.14	0.47
1:13:619:U:H3	4:3E:134:ASP:HB2	1.79	0.47
23:2L:32:G:H5''	23:2L:33:OMC:OP2	2.14	0.47
43:E8:28:SER:OG	43:E8:31:GLU:HG2	2.15	0.47
1:13:1313:U:P	19:AI:6:LYS:HD3	2.55	0.47
10:1A:30:SER:HB3	10:1A:84:GLN:OE1	2.15	0.47
26:1H:602:G:N2	26:1H:655:A:C8	2.82	0.47
31:49:64:THR:OG1	31:49:94:LEU:HD21	2.15	0.47
51:I5:42:PHE:O	51:I5:43:TYR:HB3	2.15	0.47
30:39:1:MET:HB2	30:39:2:LYS:HD3	1.96	0.47
8:72:64:LYS:O	8:72:79:VAL:HB	2.15	0.47
26:14:1200:C:H5'	58:14:3677:HOH:O	2.13	0.47
26:1H:376:C:P	58:1H:3724:HOH:O	2.73	0.47
1:1G:4:U:H3'	1:1G:5:U:H5'	1.96	0.47
26:14:959:A:N6	26:14:960:A:N1	2.62	0.47
33:69:41:GLU:HA	33:69:44:LEU:HB2	1.96	0.47
26:14:977:G:C4	26:14:978:G:C8	3.02	0.47
26:14:931:G:H3'	26:14:931:G:C8	2.49	0.47
26:14:61:G:OP2	49:G5:54:LYS:NZ	2.47	0.47
1:1G:1524:C:N4	1:1G:1525:G:O6	2.47	0.47
32:51:118:PRO:HG2	32:51:140:LYS:HE2	1.95	0.47
26:14:1167:U:O2	26:14:1183:G:N2	2.48	0.47
9:82:42:ARG:NH1	9:82:71:SER:O	2.47	0.47
51:M8:39:CYS:H	51:M8:41:PRO:HD2	1.78	0.47
2:12:6:THR:OG1	2:12:7:VAL:N	2.46	0.47
26:1H:392:C:P	58:1H:3725:HOH:O	2.72	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:1154:G:OP1	41:85:58:ARG:HD3	2.14	0.47
29:21:101:ARG:O	29:21:201:THR:OG1	2.28	0.47
26:1H:751:A:P	58:1H:3860:HOH:O	2.71	0.47
16:7I:5:ARG:HE	16:7I:22:THR:CG2	2.28	0.47
1:13:266:G:O3'	17:8I:67:LYS:HB2	2.15	0.47
26:1H:888:C:H2'	26:1H:889:C:C2	2.50	0.47
1:13:1320:C:N4	19:AI:36:ARG:HG3	2.29	0.47
5:42:7:GLU:O	5:42:34:VAL:HA	2.15	0.47
1:1G:1095:U:OP1	1:1G:1108:G:N2	2.48	0.47
2:12:84:GLU:O	2:12:219:VAL:HG21	2.15	0.47
26:14:2319:G:H4'	26:14:2320:A:O5'	2.14	0.47
1:13:1260:C:OP1	1:13:1284:C:H4'	2.14	0.47
1:1G:114:U:H2'	1:1G:115:G:H8	1.79	0.47
26:14:860:U:C2	26:14:2268:A:C8	3.02	0.47
26:1H:298:G:H5''	26:1H:299:A:OP1	2.15	0.47
26:1H:2123:G:H22	26:1H:2175:C:H42	1.62	0.47
34:58:43:THR:HB	34:58:46:VAL:HG12	1.96	0.47
1:13:136:C:O3'	16:7I:65:GLN:NE2	2.47	0.47
34:15:33:LEU:HA	34:15:38:HIS:HD2	1.80	0.47
26:1H:1969:A:H1'	26:1H:1973:G:O4'	2.14	0.47
1:1G:1129:C:C4	1:1G:1139:G:N1	2.82	0.47
1:13:622:A:C8	1:13:623:C:C6	3.03	0.47
26:14:857:C:H2'	26:14:858:U:C6	2.48	0.47
1:13:560:U:H5'	1:13:566:G:N2	2.29	0.47
26:14:2031:A:C6	26:14:2498:C:H1'	2.50	0.47
30:31:39:TRP:O	30:31:43:LYS:HG2	2.15	0.47
2:1E:59:GLU:HB2	2:1E:221:LEU:HD11	1.96	0.47
7:6E:120:ILE:O	7:6E:124:LEU:HB2	2.14	0.47
26:14:1926:U:H2'	26:14:1928:A:OP2	2.14	0.47
26:1H:270(B):A:N1	26:1H:273:G:O2'	2.44	0.47
3:22:113:ALA:HA	3:22:202:ILE:HD11	1.95	0.47
26:1H:1748:G:H2'	26:1H:1749:A:C8	2.49	0.47
26:14:298:G:OP1	45:C5:85:VAL:HA	2.15	0.47
36:35:86:LYS:HB3	36:35:118:GLY:HA3	1.97	0.47
1:13:119:A:N6	1:13:288:A:H1'	2.30	0.47
1:1G:170:U:O2'	1:1G:171:A:H5'	2.15	0.47
1:1G:1126:U:H4'	1:1G:1127:G:N7	2.29	0.47
9:8E:118:LYS:O	9:8E:119:ALA:HB3	2.15	0.47
44:B5:32:PRO:HA	44:B5:77:LYS:HB2	1.96	0.47
12:3A:54:LYS:HD2	12:3A:54:LYS:H	1.79	0.47
1:1G:518:C:H5''	1:1G:519:C:C6	2.50	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:1H:987:G:O2'	26:1H:1000:A:N3	2.41	0.47
26:14:1159:U:O2'	26:14:1160:G:H5'	2.14	0.47
24:3K:59:U:H3'	24:3K:60:U:C6	2.50	0.47
47:E5:27:GLU:HB2	47:E5:69:PHE:HD1	1.78	0.47
10:1I:77:PRO:CB	10:1I:79:ARG:HH12	2.23	0.47
31:4I:64:THR:HG22	31:4I:66:GLN:N	2.30	0.47
41:C8:95:LEU:HG	42:D8:4:ILE:HD13	1.97	0.47
50:L8:46:ASN:O	50:L8:50:VAL:HG22	2.15	0.47
2:12:54:THR:HG22	2:12:58:ILE:HD11	1.97	0.47
20:BI:61:SER:O	20:BI:65:LYS:HB2	2.15	0.47
3:22:91:LEU:HD11	3:22:101:LEU:HD12	1.97	0.47
1:13:703:G:H4'	1:13:704:A:O5'	2.15	0.47
26:14:459:U:H4'	54:L5:40:TRP:CZ3	2.49	0.47
13:4A:32:GLU:OE2	13:4A:33:ALA:N	2.47	0.47
26:1H:95:G:O2'	49:K8:48:HIS:HB3	2.15	0.47
28:19:43:ARG:HH11	28:19:43:ARG:CG	2.28	0.47
34:58:95:PRO:O	34:58:96:GLU:CD	2.53	0.47
1:1G:545:C:OP1	4:32:61:LYS:NZ	2.48	0.47
50:L8:56:VAL:HG12	50:L8:57:GLU:N	2.29	0.47
26:1H:2171:A:O2'	26:1H:2172:U:O5'	2.32	0.47
36:78:15:ARG:HA	36:78:15:ARG:HD2	1.79	0.47
1:1G:1286:A:C8	1:1G:1287:A:H4'	2.49	0.47
1:1G:1096:C:H2'	1:1G:1097:C:H6	1.80	0.47
26:14:249:C:OP1	58:14:3518:HOH:O	2.20	0.47
7:6E:71:PRO:O	7:6E:91:VAL:HG21	2.15	0.47
1:13:547:A:OP1	4:3E:73:ARG:NH2	2.48	0.47
38:55:24:GLN:HE22	38:55:36:THR:HG21	1.79	0.47
26:14:1011:G:N3	26:14:1151:G:N2	2.62	0.47
26:14:2139:C:N4	26:14:2152:G:O6	2.47	0.47
26:14:2065:C:H2'	26:14:2066:C:C6	2.50	0.47
26:1H:2244:U:H2'	26:1H:2245:U:O4'	2.15	0.47
26:1H:1705:G:O2'	26:1H:1706:U:H5'	2.14	0.47
1:13:445:G:H1	1:13:489:C:N4	2.12	0.47
10:1A:3:LYS:HD2	10:1A:77:PRO:HG3	1.97	0.47
26:14:1027:A:H8	26:14:1027:A:OP2	1.98	0.47
4:3E:108:LEU:HB3	4:3E:110:PHE:CD1	2.50	0.47
26:1H:2111:C:H5	26:1H:2147:G:C6	2.33	0.47
1:13:22:G:H2'	1:13:23:C:H6	1.80	0.47
28:11:6:PHE:CE2	28:11:18:VAL:HG23	2.48	0.47
35:68:93:PRO:HG3	35:68:114:ILE:HG12	1.96	0.47
8:72:104:ARG:O	8:72:107:LEU:HG	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1G:451:A:N6	1:1G:480:U:H2'	2.30	0.47
36:78:130:PHE:CE1	36:78:146:VAL:HG23	2.50	0.47
51:I5:48:ARG:HH12	51:I5:51:ASP:HB3	1.79	0.47
26:1H:1155:A:H5'	41:C8:55:ARG:HE	1.80	0.47
37:45:115:MET:O	37:45:119:ARG:HB2	2.15	0.47
26:14:76:C:O3'	49:G5:59:ARG:HG3	2.13	0.47
20:BA:49:ALA:O	20:BA:100:ILE:HG21	2.14	0.47
1:1G:1443:G:H22	40:75:119:LYS:HB2	1.80	0.47
3:22:120:VAL:O	3:22:123:GLN:HB3	2.14	0.47
4:3E:19:LEU:HD11	4:3E:63:LYS:HG3	1.97	0.47
26:14:2862:G:H2'	26:14:2863:C:H6	1.80	0.47
1:1G:1084:G:H5'	1:1G:1102:A:OP2	2.14	0.47
26:1H:15:G:C2	26:1H:16:G:C8	3.02	0.47
1:1G:890:G:O2'	1:1G:906:G:O6	2.29	0.47
26:1H:1690:A:H3'	26:1H:1691:C:H6	1.80	0.47
2:12:179:LYS:HA	8:72:72:PRO:HG3	1.96	0.47
1:1G:570:G:H1'	1:1G:820:U:C4	2.50	0.47
24:3L:56:C:H5'	26:14:2169:A:C8	2.49	0.47
1:13:947:G:H2'	1:13:948:C:O4'	2.14	0.47
2:1E:54:THR:HB	2:1E:201:ILE:HD11	1.97	0.47
26:14:1115:G:H2'	26:14:1116:C:C6	2.49	0.47
26:14:628:G:H5''	55:M5:18:ALA:HB2	1.96	0.47
26:1H:1753:G:N1	26:1H:1756:G:C2	2.82	0.47
1:1G:1401:G:C2	1:1G:1402:C:H1'	2.50	0.47
26:1H:185:U:H4'	26:1H:218:A:H4'	1.97	0.47
36:35:49:ARG:CZ	55:M5:58:ILE:HG22	2.45	0.47
34:58:10:GLU:HA	34:58:11:PRO:HD3	1.72	0.47
1:13:719:C:O2'	18:9I:49:LYS:HB3	2.15	0.47
26:14:2586:C:C2'	26:14:2587:A:H5'	2.44	0.47
5:4E:153:LYS:HD3	5:4E:154:GLY:H	1.79	0.47
26:1H:2475:C:H4'	26:1H:2476:A:OP1	2.13	0.47
26:14:1520:U:H2'	26:14:1521:G:O4'	2.15	0.47
9:8E:49:PRO:O	9:8E:53:VAL:HB	2.15	0.47
27:1J:10:C:C4	27:1J:11:C:C5	3.03	0.47
1:1G:568:G:N3	1:1G:574:A:H2	2.11	0.47
1:1G:142:G:H2'	1:1G:143:A:H8	1.80	0.47
37:88:41:TRP:CD1	37:88:96:VAL:HG22	2.50	0.47
48:F5:4:VAL:HG11	48:F5:11:ARG:HH12	1.80	0.47
1:1G:1446:A:OP1	1:1G:1446:A:H4'	2.14	0.47
53:O8:27:LYS:NZ	53:O8:27:LYS:H	2.13	0.47
26:1H:600:G:H2'	26:1H:601:C:O4'	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
13:4I:40:ASN:HB3	13:4I:43:THR:HG23	1.96	0.47
38:98:38:VAL:HB	38:98:39:PRO:HD3	1.97	0.47
51:I5:22:ILE:HG12	51:I5:23:GLU:H	1.80	0.47
1:1G:520:A:N1	1:1G:536:C:H1'	2.29	0.47
26:14:1022:G:C6	26:14:1140:C:C4	3.03	0.47
26:14:138:G:H22	44:B5:44:GLU:CD	2.12	0.47
55:Q8:48:PHE:CD2	55:Q8:49:VAL:HG12	2.50	0.47
22:1K:11:C:H2'	22:1K:12:U:O4'	2.14	0.47
26:1H:562:U:C4	26:1H:2036:C:O4'	2.67	0.47
8:7E:10:LEU:CD2	8:7E:83:ILE:HD11	2.40	0.47
1:1G:960:U:H4'	1:1G:961:U:H5''	1.96	0.47
19:AA:41:VAL:HB	19:AA:44:MET:HG3	1.97	0.47
30:39:123:LEU:HG	30:39:125:LEU:HD22	1.97	0.47
37:88:59:ARG:N	37:88:59:ARG:HH11	2.10	0.47
19:AA:9:VAL:HG13	51:I5:63:TYR:CE1	2.49	0.47
1:1G:1368:G:C5'	9:82:112:LYS:HB3	2.41	0.47
36:35:63:PRO:HA	55:M5:13:ARG:HA	1.95	0.47
26:14:2244:U:H2'	26:14:2245:U:O4'	2.14	0.47
34:15:15:LEU:HB2	34:15:134:ARG:CG	2.44	0.47
4:32:13:ARG:C	4:32:15:GLU:H	2.18	0.47
26:14:1669:A:H5''	26:14:1670:C:OP2	2.15	0.47
1:1G:539:A:H2'	1:1G:540:G:H8	1.80	0.47
26:14:2298:A:H61	26:14:2318:G:H2'	1.80	0.47
31:49:121:ASN:OD1	31:49:124:SER:N	2.46	0.47
31:41:128:ARG:NH2	31:41:128:ARG:HB2	2.30	0.47
39:A8:15:ARG:HD2	39:A8:88:ASP:OD2	2.14	0.47
1:1G:624:C:H2'	1:1G:625:G:H8	1.80	0.47
1:13:598:U:H2'	1:13:599:C:H6	1.80	0.47
53:K5:27:LYS:NZ	53:K5:27:LYS:HB3	2.29	0.47
30:39:107:LYS:HE2	30:39:205:ARG:HD2	1.97	0.47
20:BI:14:LYS:HG3	20:BI:17:ARG:HH21	1.79	0.47
39:65:69:VAL:O	39:65:72:ALA:HB3	2.15	0.47
1:13:1455:G:OP1	20:BI:35:THR:OG1	2.11	0.47
12:3A:49:ASN:N	12:3A:49:ASN:OD1	2.48	0.47
46:D5:108:PRO:HG3	46:D5:142:SER:HB2	1.97	0.47
26:1H:1127:A:N7	26:1H:2488:A:O2'	2.43	0.47
26:14:389:G:H22	36:35:72:PRO:HD3	1.79	0.47
1:1G:685:G:N1	1:1G:686:U:O4	2.47	0.47
1:13:971:G:N2	1:13:1363:A:OP2	2.45	0.47
33:61:95:LYS:HE3	33:61:99:GLU:OE1	2.15	0.47
27:1J:97:G:H2'	27:1J:98:G:O4'	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:57:C:H2'	26:14:58:G:O4'	2.15	0.47
26:1H:278:A:H5''	26:1H:278:A:H8	1.79	0.47
1:13:683:G:H2'	1:13:684:A:C8	2.50	0.47
45:G8:89:PHE:HD1	45:G8:90:LEU:N	2.13	0.47
48:F5:82:LEU:H	48:F5:82:LEU:HD23	1.79	0.47
23:2L:38:A:H2'	23:2L:39:A:O4'	2.15	0.47
1:13:741:G:H2'	1:13:742:G:O4'	2.15	0.47
12:3A:54:LYS:HD2	12:3A:54:LYS:N	2.29	0.47
1:1G:1324:A:H4'	1:1G:1362:C:H4'	1.96	0.47
44:F8:3:THR:CB	44:F8:6:ASP:HB2	2.45	0.47
31:41:64:THR:HG21	31:41:66:GLN:HB2	1.97	0.47
26:1H:1774:C:H5''	58:1H:3750:HOH:O	2.15	0.47
49:K8:15:LYS:H	49:K8:67:LYS:HZ1	1.56	0.47
26:14:276:A:N3	26:14:277:C:N4	2.63	0.47
27:1J:76:G:H2'	27:1J:77:U:O4'	2.15	0.47
55:M5:14:VAL:HG11	55:M5:22:VAL:HG13	1.97	0.47
43:E8:12:ILE:HG13	43:E8:42:ARG:NH1	2.27	0.47
28:19:71:ASP:OD1	28:19:71:ASP:N	2.47	0.47
26:1H:1657:C:H2'	26:1H:1658:C:H6	1.79	0.47
4:32:4:TYR:HE2	4:32:11:LEU:HD21	1.80	0.47
12:3I:89:ARG:HG2	12:3I:90:VAL:N	2.30	0.47
26:14:2286:A:C8	26:14:2287:A:N6	2.83	0.47
19:AA:9:VAL:HG13	51:I5:63:TYR:HE1	1.80	0.47
26:14:779:U:OP1	28:19:49:ILE:HG22	2.15	0.47
26:1H:241:A:H4'	26:1H:242:G:OP1	2.14	0.47
26:14:2361:A:OP1	55:M5:27:THR:HG23	2.14	0.47
26:1H:1324:G:N2	26:1H:1331:A:C4	2.83	0.47
26:1H:2164:C:H41	26:1H:2165:G:N2	2.13	0.47
1:13:639:G:C2	1:13:640:A:C5	3.03	0.47
24:3L:52:G:H1	24:3L:62:C:N4	2.13	0.47
26:14:818:G:H4'	26:14:838:C:O3'	2.15	0.47
11:2I:112:THR:O	11:2I:114:VAL:HG12	2.14	0.47
26:1H:768:G:O2'	26:1H:1379:A:N6	2.48	0.47
26:1H:2074:U:OP1	58:1H:3681:HOH:O	2.20	0.47
36:78:96:THR:HG22	36:78:126:VAL:HG21	1.97	0.47
28:19:181:GLU:HA	28:19:272:ALA:HB1	1.97	0.47
7:6E:5:ARG:CZ	7:6E:7:ALA:HA	2.45	0.47
1:1G:547:A:OP2	4:32:2:GLY:N	2.48	0.47
34:15:91:LEU:O	34:15:95:PRO:HB3	2.14	0.47
26:14:1167:U:C2	26:14:1183:G:N2	2.83	0.47
1:13:684:A:C6	1:13:685:G:C6	3.03	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:13:1074:G:O2'	1:13:1101:A:N1	2.42	0.47
12:3A:109:GLY:HA3	12:3A:121:GLY:O	2.15	0.47
26:1H:181:A:H1'	26:1H:435:C:H5'	1.96	0.47
26:1H:2774:C:H2'	26:1H:2775:A:O4'	2.15	0.47
8:7E:82:HIS:CE1	8:7E:136:GLU:OE2	2.68	0.47
1:13:501:C:H2'	1:13:502:G:H8	1.80	0.47
1:13:509:A:H5''	4:3E:55:ALA:HB2	1.97	0.47
34:15:103:VAL:HG11	34:15:120:LEU:HD22	1.96	0.47
40:75:27:THR:HG23	40:75:89:VAL:HG22	1.96	0.47
54:P8:35:ARG:NH1	54:P8:42:LEU:HD11	2.30	0.47
1:13:321:A:C2	1:13:333:G:C2	3.03	0.47
29:21:33:VAL:CG1	29:21:89:ASP:HA	2.45	0.47
15:6I:10:LYS:HD2	15:6I:10:LYS:HA	1.68	0.47
26:14:654:A:OP1	26:14:654:A:H8	1.98	0.47
24:1L:58:A:O2'	24:1L:61:C:N4	2.48	0.47
33:61:86:THR:HA	33:61:123:LEU:HD13	1.96	0.47
2:12:48:MET:HA	2:12:51:LEU:HB2	1.97	0.47
1:1G:718:G:H5'	11:2A:117:ASN:ND2	2.29	0.47
26:14:1576:U:H2'	26:14:1577:C:H6	1.79	0.47
31:41:112:PRO:HA	51:M8:37:SER:HB2	1.97	0.47
2:12:213:LEU:HA	2:12:216:SER:HB3	1.96	0.47
32:59:88:LEU:HB2	32:59:163:TYR:HB2	1.95	0.47
24:3K:72:C:H3'	24:3K:73:A:C5'	2.37	0.47
26:1H:2598:A:P	58:1H:3656:HOH:O	2.69	0.47
24:3K:18:G:H1'	24:3K:58:A:C2	2.50	0.47
26:1H:1515:C:H2'	26:1H:1516:U:C6	2.49	0.47
39:65:78:LEU:HD21	39:65:108:GLY:HA3	1.97	0.47
55:M5:6:THR:HG22	55:M5:59:LYS:O	2.15	0.47
1:13:765:G:H5''	1:13:766:A:OP1	2.15	0.47
26:1H:1533:C:C2	26:1H:1534:G:N2	2.83	0.47
26:1H:598:G:H1	26:1H:659:C:H42	1.61	0.47
37:88:37:LEU:HD21	37:88:130:LYS:HE3	1.96	0.47
1:1G:1187:G:H4'	9:82:111:ARG:HH11	1.79	0.47
39:65:14:VAL:HG21	39:65:90:GLY:O	2.15	0.47
1:1G:1258:G:H2'	1:1G:1259:C:H6	1.80	0.47
1:13:342:C:H42	1:13:347:G:H1	1.63	0.47
26:14:654(A):A:H2'	26:14:654(B):C:C6	2.50	0.47
47:E5:25:ARG:HD2	47:E5:29:GLN:NE2	2.29	0.47
3:2E:6:HIS:HE1	3:2E:184:TYR:CD2	2.33	0.47
3:22:21:ARG:HH11	3:22:21:ARG:HB3	1.80	0.47
20:BI:50:GLU:HG3	20:BI:100:ILE:HG12	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:1E:97:TRP:HZ2	2:1E:102:LEU:HD13	1.79	0.47
35:25:68:GLU:OE2	35:25:78:ARG:NH1	2.47	0.47
23:2K:8:4SU:OP2	23:2K:8:4SU:H6	2.14	0.47
1:1G:560:U:H5'	1:1G:566:G:N2	2.29	0.47
24:3L:62:C:O5'	24:3L:62:C:H6	1.97	0.47
26:1H:1337:G:C4	26:1H:1338:G:C8	3.03	0.47
26:1H:989:G:C8	50:L8:13:ILE:HD11	2.49	0.47
26:1H:2788:C:P	29:21:61:ARG:HH22	2.37	0.47
26:1H:248:G:H5'	26:1H:250:G:N7	2.30	0.47
26:14:2228:G:C6	26:14:2229:C:C4	3.03	0.47
26:1H:942:G:OP2	36:78:39:LYS:HE2	2.14	0.47
26:14:2808:U:H2'	26:14:2809:A:H8	1.79	0.47
5:42:147:ASP:O	5:42:151:LEU:HG	2.14	0.47
26:14:2757:A:C6	32:59:67:LEU:HD22	2.50	0.47
20:BA:69:GLY:O	20:BA:73:HIS:NE2	2.48	0.47
27:16:28:C:H2'	27:16:29:A:H8	1.79	0.47
36:78:114:ILE:HG12	36:78:130:PHE:CD2	2.50	0.47
1:13:1442:G:H2'	1:13:1443:G:H5'	1.97	0.47
50:H5:6:VAL:O	50:H5:34:GLU:HA	2.15	0.47
26:14:1779:U:H5''	26:14:1780:A:C5'	2.45	0.47
32:59:120:GLY:O	32:59:135:GLY:HA2	2.15	0.47
1:1G:176:C:O2'	1:1G:177:C:H5'	2.16	0.47
38:98:70:LEU:HD23	38:98:70:LEU:HA	1.72	0.47
2:12:30:ARG:NH1	2:12:194:PRO:HB2	2.30	0.47
26:1H:1093:G:H1'	26:1H:1099:G:H1	1.80	0.47
26:14:1788:C:H2'	26:14:1789:A:H8	1.79	0.47
30:39:153:SER:HB2	30:39:190:GLU:H	1.79	0.47
26:14:2176:A:H2'	26:14:2177:C:H6	1.79	0.47
26:1H:59:U:O2'	26:1H:73:A:H2'	2.15	0.47
26:14:372:G:OP2	48:F5:69:LYS:NZ	2.46	0.47
26:14:1449(A):G:H2'	26:14:1450:C:H6	1.80	0.47
5:4E:43:LEU:HD23	5:4E:133:TYR:CD1	2.50	0.47
24:3L:71:G:H2'	24:3L:72:C:C5'	2.45	0.46
1:13:1367:C:N3	1:13:1368:G:C8	2.84	0.46
1:13:1392:G:N2	1:13:1502:A:H8	2.14	0.46
26:14:1043:C:H2'	26:14:1044:G:H5'	1.98	0.46
5:42:103:GLY:C	5:42:106:PRO:HD2	2.35	0.46
26:14:1225:C:H4'	42:95:85:LYS:CB	2.45	0.46
10:1I:77:PRO:HB2	10:1I:79:ARG:NH1	2.23	0.46
27:1J:24:G:C8	27:1J:56:G:C5	3.03	0.46
24:3K:6:G:H22	24:3K:67:C:H42	1.63	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
12:3I:126:LYS:HG3	12:3I:128:ALA:N	2.26	0.46
34:15:35:ARG:HB2	34:15:42:TRP:HZ3	1.79	0.46
26:1H:1142:U:H5'	26:1H:1142(A):A:H8	1.79	0.46
26:14:1071:G:H1'	26:14:1089:G:H2'	1.96	0.46
26:14:270(I):G:H2'	26:14:270(J):G:C8	2.43	0.46
26:14:2327:A:H2'	26:14:2328:A:H8	1.79	0.46
34:58:96:GLU:C	34:58:98:VAL:N	2.68	0.46
37:88:106:VAL:HG21	37:88:114:ALA:CB	2.44	0.46
1:13:344:A:O2'	1:13:346:G:N7	2.47	0.46
32:59:121:ILE:HD12	32:59:133:VAL:HG13	1.97	0.46
26:14:528:A:C2	26:14:2043:C:H4'	2.50	0.46
26:1H:639:U:O2'	26:1H:640:C:H5'	2.14	0.46
24:3L:8:U:H4'	24:3L:9:A:OP1	2.14	0.46
28:19:37:LEU:CB	28:19:38:LYS:HG2	2.45	0.46
1:13:626:U:C2	1:13:627:G:C8	3.03	0.46
1:13:1190:G:OP2	3:2E:5:ILE:HG23	2.15	0.46
13:4A:92:HIS:NE2	13:4A:98:VAL:HG21	2.30	0.46
45:C5:50:ARG:HG2	45:C5:53:PRO:HG3	1.97	0.46
16:7I:67:THR:HG22	16:7I:68:ASP:H	1.80	0.46
26:14:1003:G:O2'	26:14:1010:A:N1	2.34	0.46
4:3E:8:VAL:CG1	4:3E:21:LEU:HB2	2.45	0.46
20:BA:19:SER:O	20:BA:23:ARG:N	2.43	0.46
2:1E:60:ASP:HB3	2:1E:64:ARG:NH1	2.30	0.46
1:13:939:G:C6	1:13:940:C:N4	2.83	0.46
4:3E:190:ASP:HB3	4:3E:192:GLU:OE1	2.15	0.46
42:95:22:VAL:HG22	42:95:23:GLU:H	1.80	0.46
20:BA:53:LEU:O	20:BA:57:ARG:HG3	2.14	0.46
40:B8:107:ASP:OD2	40:B8:109:GLU:HG3	2.14	0.46
1:1G:408:A:H2'	1:1G:409:G:O4'	2.16	0.46
16:7A:57:ARG:HA	16:7A:60:LEU:HD12	1.96	0.46
4:32:14:ARG:HG3	4:32:14:ARG:HH11	1.80	0.46
38:55:55:ALA:HA	38:55:80:PHE:CZ	2.50	0.46
46:D5:10:ARG:HH21	46:D5:26:GLY:H	1.61	0.46
9:82:63:ILE:HD11	9:82:81:ILE:HD11	1.96	0.46
26:1H:945:A:OP2	26:1H:945:A:H4'	2.15	0.46
24:3K:2:C:N3	24:3K:72:C:N4	2.62	0.46
40:75:64:ARG:NH1	40:75:103:ARG:HA	2.28	0.46
42:95:71:LEU:HA	42:95:86:GLY:HA2	1.97	0.46
29:29:111:ARG:HD2	29:29:160:TYR:CD2	2.51	0.46
1:1G:837:G:H1	1:1G:849:C:N4	2.12	0.46
26:1H:1064:C:N4	26:1H:1070:A:OP1	2.48	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
33:61:21:VAL:CG2	33:61:25:TYR:HD2	2.29	0.46
26:14:1770:G:H1	26:14:1982:C:H42	1.63	0.46
2:12:165:VAL:O	2:12:167:PRO:HD3	2.15	0.46
1:1G:1299:A:N1	1:1G:1301:U:N3	2.64	0.46
30:31:129:PHE:O	30:31:130:ALA:HB3	2.16	0.46
26:1H:527:C:H4'	26:1H:528:A:H5'	1.96	0.46
5:4E:75:THR:OG1	5:4E:76:ILE:N	2.41	0.46
24:1L:10:G:O2'	24:1L:11:C:OP1	2.31	0.46
41:C8:108:GLU:OE1	41:C8:112:ARG:NH1	2.48	0.46
26:14:1006:C:O2	34:15:106:MET:HB3	2.15	0.46
26:14:839:U:H2'	26:14:840:C:C6	2.51	0.46
27:16:31:C:H2'	27:16:32:C:H6	1.79	0.46
2:1E:215:LEU:O	2:1E:219:VAL:HG23	2.15	0.46
26:14:1716:U:H2'	26:14:1717:G:C8	2.50	0.46
1:1G:442:C:H2'	1:1G:443:C:C6	2.50	0.46
26:1H:2707:G:O3'	38:98:68:ARG:HG2	2.15	0.46
12:3I:71:PRO:O	12:3I:102:ARG:HD3	2.15	0.46
26:14:925:C:H2'	26:14:926:A:H8	1.79	0.46
26:14:107:C:H2'	26:14:108:U:H6	1.80	0.46
26:14:2577:A:H2'	26:14:2614:A:N6	2.31	0.46
1:13:1527:C:N4	1:13:1528:U:O4	2.48	0.46
26:1H:775:G:C5	26:1H:794:G:C8	3.03	0.46
26:14:510:C:H2'	26:14:511:U:O4'	2.15	0.46
1:1G:149:A:H2'	1:1G:150:C:C6	2.50	0.46
28:11:174:ILE:HD12	28:11:174:ILE:N	2.30	0.46
29:21:14:ILE:HA	29:21:14:ILE:HD12	1.58	0.46
1:1G:340:U:H2'	1:1G:341:C:C6	2.49	0.46
26:14:903:C:H2'	26:14:904:C:C6	2.50	0.46
18:9A:34:TYR:HA	18:9A:40:LEU:HD11	1.97	0.46
36:78:29:LYS:HD2	36:78:30:THR:HG22	1.98	0.46
55:Q8:26:LYS:HD3	55:Q8:26:LYS:O	2.15	0.46
39:A8:34:HIS:CD2	39:A8:54:LEU:HD23	2.50	0.46
26:14:1154:G:OP2	41:85:58:ARG:NH1	2.48	0.46
27:1J:44:G:C2	27:1J:48:A:C2	3.03	0.46
42:95:76:LYS:NZ	42:95:82:ARG:HE	2.14	0.46
1:13:57:G:H2'	1:13:58:C:C6	2.50	0.46
26:1H:1534:G:N2	26:1H:1538:G:H22	2.11	0.46
37:88:51:ARG:NH1	37:88:52:VAL:HG23	2.25	0.46
32:51:4:ILE:HG13	32:51:6:ARG:HB2	1.96	0.46
26:1H:963:U:OP1	58:1H:3797:HOH:O	2.21	0.46
26:1H:459:U:H4'	54:P8:40:TRP:CZ3	2.50	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
38:98:10:LEU:O	38:98:12:ARG:N	2.49	0.46
5:42:83:GLU:HB3	5:42:88:LYS:HG3	1.98	0.46
1:13:447:G:O5'	1:13:447:G:H8	1.99	0.46
1:13:413:G:N2	1:13:428:G:H1'	2.29	0.46
46:D5:128:VAL:HG23	46:D5:160:GLY:HA3	1.97	0.46
32:51:154:PRO:HB3	32:51:163:TYR:CE2	2.49	0.46
34:58:35:ARG:O	34:58:42:TRP:HZ3	1.98	0.46
32:59:76:VAL:HA	32:59:79:VAL:HG22	1.97	0.46
7:6E:5:ARG:NH1	7:6E:8:GLU:OE2	2.48	0.46
26:1H:1789:A:H5'	28:11:220:HIS:O	2.16	0.46
23:2L:4:G:H1	23:2L:70:C:H42	1.63	0.46
26:1H:1068:G:H1'	26:1H:1096:A:N3	2.30	0.46
4:3E:64:LEU:HD22	4:3E:198:VAL:HG11	1.96	0.46
26:1H:1567:A:H5'	28:11:58:HIS:ND1	2.30	0.46
6:5E:23:LYS:HA	6:5E:26:ILE:HD12	1.97	0.46
33:61:113:ARG:H	33:61:116:LEU:HD13	1.80	0.46
54:L5:6:GLN:HA	54:L5:7:PRO:HD3	1.72	0.46
40:75:18:ASP:OD1	40:75:19:LEU:HG	2.15	0.46
26:1H:2101:G:H1	26:1H:2188:C:H42	1.63	0.46
1:1G:1449:C:H3'	1:1G:1450:U:H4'	1.97	0.46
2:12:9:GLU:HB2	2:12:217:ARG:NH2	2.30	0.46
1:1G:536:C:H2'	1:1G:537:G:C8	2.51	0.46
26:1H:1664:A:OP1	58:1H:3715:HOH:O	2.19	0.46
26:1H:2061:G:C2	26:1H:2063:C:C4	3.03	0.46
26:1H:2062:A:N6	26:1H:2503:A:H62	2.13	0.46
36:78:138:LEU:CD1	36:78:144:GLU:HG3	2.40	0.46
1:1G:617:G:C2	1:1G:618:C:C5	3.04	0.46
26:14:739:G:P	58:14:3506:HOH:O	2.70	0.46
1:1G:1352:C:H2'	1:1G:1353:G:C8	2.51	0.46
34:15:39:ARG:HA	34:15:40:PRO:HD2	1.70	0.46
26:1H:1142(A):A:H4'	34:58:25:ARG:NH2	2.29	0.46
26:14:1087:G:H2'	26:14:1089:G:H1'	1.97	0.46
16:7I:26:ARG:HH21	16:7I:31:LYS:HD3	1.80	0.46
26:1H:960:A:C8	26:1H:962:G:C8	3.02	0.46
1:1G:1372:U:H2'	1:1G:1373:G:O4'	2.15	0.46
1:13:1063:C:H3'	1:13:1064:G:H2'	1.97	0.46
26:14:96:G:H4'	49:G5:48:HIS:ND1	2.30	0.46
36:78:1:MET:HE1	36:78:6:LEU:HA	1.96	0.46
1:13:954:G:H2'	1:13:955:U:C6	2.50	0.46
6:52:97:PHE:HE2	18:9A:62:GLU:HG2	1.81	0.46
26:1H:581:C:OP1	41:C8:33:ARG:HG3	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:1H:2457:U:C2'	26:1H:2458:G:H5'	2.45	0.46
1:1G:1329:A:H5'	13:4A:29:ARG:HE	1.80	0.46
31:49:56:ALA:HB2	31:49:153:ARG:CZ	2.46	0.46
23:2L:20:G:OP1	23:2L:61:U:N3	2.46	0.46
28:11:17:THR:CG2	28:11:205:VAL:H	2.28	0.46
49:K8:28:LYS:HD2	49:K8:56:GLN:NE2	2.30	0.46
26:14:656:G:H2'	26:14:657:U:O4'	2.15	0.46
26:1H:297:C:H2'	26:1H:298:G:O4'	2.16	0.46
1:1G:448:A:OP2	1:1G:485:G:N2	2.40	0.46
26:1H:2107:C:O2	26:1H:2182:G:N2	2.29	0.46
11:2I:112:THR:HA	11:2I:113:PRO:HD3	1.68	0.46
47:E5:12:ASN:HA	47:E5:14:ARG:NH2	2.31	0.46
1:13:475:G:H2'	1:13:476:G:O4'	2.16	0.46
19:AI:6:LYS:HE2	19:AI:6:LYS:HB3	1.77	0.46
27:1J:99:A:C5	27:1J:100:G:C8	3.03	0.46
26:1H:1519:G:C2'	26:1H:1520:U:H5'	2.45	0.46
18:9I:32:ARG:HH11	18:9I:65:ILE:HD13	1.81	0.46
1:1G:37:U:O2'	1:1G:547:A:N1	2.43	0.46
1:1G:801:U:H2'	1:1G:802:A:H8	1.80	0.46
48:F5:4:VAL:HG11	48:F5:11:ARG:NH1	2.30	0.46
1:13:1292:U:H2'	1:13:1293:G:C8	2.50	0.46
26:1H:483:A:O4'	45:G8:48:ALA:HB1	2.16	0.46
19:AI:52:TYR:HA	19:AI:56:GLN:O	2.15	0.46
1:13:853:G:H5''	1:13:853:G:H8	1.80	0.46
46:H8:5:LEU:HD23	46:H8:47:VAL:HG21	1.96	0.46
26:14:1166:C:H1'	26:14:1184:G:N2	2.30	0.46
42:95:1:MET:HB3	42:95:42:GLY:HA3	1.96	0.46
26:1H:2748:A:C2	32:51:63:SER:HB3	2.50	0.46
34:58:137:LYS:HE3	34:58:138:LEU:HB2	1.98	0.46
33:61:8:PRO:HA	33:61:14:ASP:HA	1.96	0.46
29:29:31:CYS:O	29:29:91:VAL:HG22	2.15	0.46
47:I8:60:PHE:CD1	47:I8:60:PHE:N	2.83	0.46
2:12:7:VAL:HG13	2:12:8:LYS:HG3	1.98	0.46
1:1G:516:U:O2'	1:1G:519:C:N3	2.49	0.46
36:78:62:LEU:HB3	55:Q8:23:VAL:HG21	1.96	0.46
26:14:1970:A:P	58:14:3577:HOH:O	2.70	0.46
22:1K:23:A:H2'	22:1K:24:G:O4'	2.14	0.46
11:2A:54:ARG:NH1	24:3L:39:U:O3'	2.48	0.46
39:A8:87:PHE:HZ	39:A8:98:VAL:HG12	1.81	0.46
52:N8:38:ALA:HB3	52:N8:40:LYS:HE3	1.96	0.46
1:1G:728:A:N7	15:6A:54:ARG:HD2	2.31	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
29:21:105:THR:HG21	29:21:164:ARG:CZ	2.46	0.46
12:3I:91:LYS:O	12:3I:91:LYS:HG3	2.14	0.46
26:1H:1171:G:N2	26:1H:1178:C:N3	2.63	0.46
46:D5:126:VAL:HG12	46:D5:163:LEU:HB3	1.98	0.46
26:14:2295:C:H5	39:65:13:ARG:NH2	2.14	0.46
1:1G:1297:C:OP1	13:4A:13:LYS:HG3	2.15	0.46
13:4A:3:ARG:HD3	13:4A:7:VAL:HA	1.97	0.46
29:21:37:ARG:O	29:21:45:THR:HA	2.16	0.46
26:1H:2314:C:H5'	31:41:38:VAL:HG21	1.96	0.46
1:1G:390:C:H2'	1:1G:391:G:C8	2.51	0.46
37:45:86:GLY:O	37:45:88:GLY:N	2.47	0.46
1:1G:1226:C:H4'	19:AA:80:TYR:CZ	2.51	0.46
4:3E:108:LEU:HD21	4:3E:183:GLY:HA3	1.96	0.46
4:3E:106:TYR:HE2	4:3E:107:ARG:NH1	2.14	0.46
26:1H:1443:G:N2	26:1H:1549:C:C2	2.84	0.46
26:1H:2863:C:H2'	26:1H:2864:G:C8	2.50	0.46
1:13:1364:U:O2'	1:13:1365:G:H5'	2.15	0.46
26:14:2867:G:OP2	40:75:119:LYS:NZ	2.34	0.46
1:1G:403:C:H42	1:1G:547:A:H5'	1.79	0.46
34:15:90:MET:HB3	34:15:98:VAL:HG12	1.98	0.46
26:1H:1530:G:O6	26:1H:1542:G:N2	2.48	0.46
26:1H:600:G:N2	26:1H:605:C:O3'	2.48	0.46
26:1H:1400:G:H2'	26:1H:1401:G:C8	2.51	0.46
11:2A:109:VAL:HG22	18:9A:86:VAL:HG13	1.97	0.46
16:7A:15:PRO:O	16:7A:16:HIS:HD2	1.99	0.46
31:49:6:ALA:O	31:49:10:LYS:N	2.38	0.46
26:1H:556:G:H2'	26:1H:557:U:C6	2.50	0.46
11:2I:29:ILE:HG13	11:2I:44:SER:HB3	1.97	0.46
4:32:126:ILE:HG22	4:32:127:THR:H	1.79	0.46
26:14:239:U:H2'	26:14:240:G:O4'	2.15	0.46
23:2K:23:G:H2'	23:2K:24:C:H6	1.81	0.46
23:2K:26:C:H2'	23:2K:27:G:O4'	2.15	0.46
39:65:15:ARG:O	39:65:19:LYS:HD3	2.14	0.46
26:14:2547:U:O2'	26:14:2548:G:H5'	2.16	0.46
36:78:76:LYS:HA	36:78:76:LYS:HD3	1.57	0.46
45:G8:34:LYS:HG2	45:G8:34:LYS:O	2.16	0.46
32:59:27:LYS:HA	32:59:27:LYS:HD2	1.69	0.46
41:C8:90:VAL:HG22	42:D8:39:LEU:HB3	1.98	0.46
44:F8:34:ALA:HA	44:F8:38:GLU:OE1	2.16	0.46
26:14:623:G:H2'	26:14:624:C:C6	2.50	0.46
11:2I:78:GLN:O	11:2I:103:LEU:HA	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:82:17:VAL:HG22	9:82:63:ILE:HD11	1.96	0.46
45:G8:94:LYS:HD2	45:G8:94:LYS:HA	1.68	0.46
22:1K:76:A:H1'	26:1H:2583:G:N2	2.30	0.46
45:C5:39:VAL:HG23	45:C5:41:GLY:N	2.24	0.46
28:11:69:ARG:HD3	28:11:105:ILE:HD11	1.97	0.46
26:14:276:A:H2'	26:14:277:C:C5	2.50	0.46
1:1G:1369:C:H2'	1:1G:1370:G:O4'	2.16	0.46
1:1G:707:C:H2'	1:1G:708:C:H6	1.81	0.46
17:8I:67:LYS:O	17:8I:68:ARG:HB3	2.16	0.46
26:14:2388:A:C2'	26:14:2389:G:H5'	2.45	0.46
29:29:33:VAL:HG11	29:29:36:ARG:NH2	2.30	0.46
29:29:87:GLU:O	29:29:89:ASP:HA	2.15	0.46
26:14:2394:C:H2'	26:14:2395:C:H6	1.79	0.46
26:1H:1278:A:H5''	38:98:36:THR:HG22	1.98	0.46
46:D5:71:VAL:HB	46:D5:88:PHE:CE1	2.51	0.46
13:4I:23:TYR:HE2	13:4I:71:ARG:HG3	1.79	0.46
38:55:29:LEU:HD12	38:55:29:LEU:HA	1.81	0.46
3:22:134:ILE:HD13	3:22:134:ILE:HA	1.81	0.46
8:72:11:THR:HG23	8:72:14:ARG:HH12	1.80	0.46
1:13:221:C:H2'	1:13:222:U:C6	2.49	0.46
29:29:26:ILE:HG21	29:29:28:ALA:HB2	1.96	0.46
49:K8:28:LYS:HD3	49:K8:53:LEU:HD21	1.98	0.46
24:3L:52:G:H1	24:3L:62:C:H42	1.63	0.46
26:1H:234:C:C2	26:1H:235:U:C5	3.03	0.46
1:1G:358:U:H2'	1:1G:359:U:H6	1.80	0.46
26:1H:2182:G:H2'	26:1H:2183:C:C6	2.50	0.46
5:42:122:GLU:OE1	5:42:131:ILE:HG13	2.16	0.46
55:M5:29:LYS:HB3	55:M5:44:LYS:HB2	1.98	0.46
17:8I:88:TYR:CD2	17:8I:89:LEU:HD22	2.49	0.46
1:1G:1453:G:O6	20:BA:51:GLU:HB3	2.16	0.46
20:BA:75:ASN:HA	20:BA:78:ALA:HB3	1.97	0.46
1:1G:1111:A:H2'	1:1G:1112:C:C6	2.51	0.46
26:1H:487:C:C2'	26:1H:488:G:H5'	2.46	0.46
34:15:34:LEU:HD21	34:15:120:LEU:HB2	1.98	0.46
1:13:834:C:C2	1:13:853:G:C2	3.04	0.46
10:1A:6:ILE:HG22	10:1A:98:ILE:HG12	1.98	0.46
41:85:19:LYS:O	41:85:22:LYS:HG3	2.16	0.46
3:22:61:ALA:C	3:22:63:ASN:H	2.19	0.46
26:14:695:G:C2	26:14:768:G:C5	3.04	0.46
2:1E:130:ARG:H	2:1E:130:ARG:HG2	1.50	0.46
24:1L:22:G:H8	24:1L:22:G:OP2	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
28:19:146:GLU:HB2	28:19:189:CYS:HB3	1.95	0.46
26:14:1665:A:H4'	35:25:67:LYS:HB2	1.98	0.46
26:1H:118:A:H3'	26:1H:119:A:H5''	1.97	0.46
26:1H:1952:A:C6	35:68:22:ILE:HD12	2.50	0.46
30:39:165:ARG:HA	30:39:168:ARG:HD3	1.97	0.46
26:14:895:U:H4'	26:14:896:A:C5	2.51	0.46
1:1G:1418:A:H2	26:14:1948:G:N3	2.13	0.46
34:58:65:LYS:HE3	34:58:65:LYS:HB2	1.56	0.46
26:14:1805:U:C2	26:14:1806:C:C5	3.03	0.46
30:39:36:VAL:HG11	30:39:183:VAL:HG11	1.96	0.46
1:13:297:G:H4'	1:13:557:G:H4'	1.98	0.46
28:19:51:VAL:HG12	28:19:54:ARG:HD2	1.97	0.46
51:M8:37:SER:OG	51:M8:42:PHE:HB3	2.16	0.46
26:1H:312:G:H5'	26:1H:331:A:O2'	2.15	0.46
46:H8:125:LEU:HG	46:H8:164:ALA:HB3	1.98	0.46
26:1H:58:G:N2	26:1H:70:G:C5	2.84	0.46
26:1H:576:U:H5	58:1H:3921:HOH:O	1.98	0.46
1:13:60:A:H4'	1:13:61:G:H5'	1.98	0.46
1:13:1177:G:C8	1:13:1178:G:C2	3.03	0.46
26:1H:1317:A:H2'	26:1H:1318:C:C6	2.51	0.46
26:1H:2376:A:N1	39:A8:87:PHE:HB3	2.31	0.46
2:12:42:ILE:HD13	2:12:43:ASP:N	2.31	0.46
26:14:1435:G:H1	26:14:1557:C:H42	1.64	0.46
13:4I:3:ARG:HB2	51:M8:34:GLU:OE2	2.15	0.46
29:21:119:ARG:HB3	29:21:120:TRP:CD1	2.51	0.46
1:13:632:A:H8	1:13:633:G:N9	2.14	0.46
26:1H:1141:U:H6	34:58:63:THR:OG1	1.99	0.46
1:13:79:G:O2'	1:13:90:C:O2	2.25	0.46
26:14:2306:C:H3'	26:14:2307:G:H5''	1.98	0.46
31:49:16:ARG:O	31:49:20:ILE:HG13	2.15	0.46
1:13:991:U:C4	1:13:1212:U:H1'	2.51	0.46
33:69:125:GLU:OE1	33:69:141:LYS:HG3	2.16	0.46
2:12:22:LYS:C	2:12:24:TRP:H	2.17	0.46
26:1H:2849:U:O4	40:B8:23:ARG:NH2	2.48	0.46
9:8E:46:ALA:HA	9:8E:78:LYS:HB2	1.97	0.46
26:1H:2012:G:OP1	43:E8:11:ARG:NH2	2.43	0.46
28:19:37:LEU:HA	28:19:38:LYS:HG2	1.98	0.46
1:13:280:C:C2	17:8I:38:ARG:HG3	2.51	0.46
26:1H:2336:A:N6	47:I8:43:THR:HB	2.31	0.46
26:14:817:C:O2'	26:14:839:U:H5''	2.16	0.46
1:1G:922:G:H4'	5:42:20:GLN:HA	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:1H:1058:U:H3	26:1H:1080:A:N6	2.13	0.46
29:29:41:LYS:HG3	29:29:42:ASP:H	1.81	0.46
1:1G:1194:U:H2'	1:1G:1195:C:H6	1.79	0.46
8:7E:58:TYR:O	8:7E:59:LEU:HD23	2.15	0.46
51:M8:49:PHE:CD2	51:M8:50:VAL:HG12	2.50	0.46
26:14:1569:A:H5'	28:19:61:LEU:CD2	2.46	0.46
26:14:2469:A:C2	26:14:2470:G:C5	3.04	0.46
1:13:319:G:H2'	1:13:320:C:O4'	2.16	0.46
1:13:37:U:O2'	1:13:500:G:H4'	2.15	0.46
46:D5:76:LEU:H	46:D5:76:LEU:HD23	1.81	0.46
46:D5:76:LEU:HA	46:D5:83:PRO:HA	1.98	0.46
4:3E:99:SER:O	4:3E:140:VAL:HG22	2.16	0.46
2:12:189:ASP:N	2:12:189:ASP:OD1	2.49	0.46
36:35:98:GLU:HA	36:35:101:VAL:HG12	1.98	0.46
9:82:17:VAL:HA	9:82:63:ILE:HG12	1.98	0.46
1:1G:590:C:H2'	1:1G:591:U:H6	1.81	0.46
1:13:452:A:H2'	1:13:453:A:C8	2.51	0.46
26:1H:1665:A:N6	58:1H:3964:HOH:O	2.48	0.46
26:14:307:G:N2	26:14:310:A:OP2	2.47	0.46
26:1H:1317:A:H2'	26:1H:1318:C:H6	1.81	0.46
26:1H:2210:G:H5''	26:1H:2211:G:N7	2.31	0.46
26:14:2688:U:H1'	26:14:2721:A:N6	2.31	0.46
30:31:125:LEU:HD21	30:31:199:TRP:CE3	2.51	0.46
9:8E:3:GLN:OE1	9:8E:20:ARG:NH1	2.48	0.46
1:1G:1065:U:C5	1:1G:1190:G:H1'	2.50	0.46
48:F5:92:LYS:O	48:F5:94:LEU:N	2.48	0.46
26:1H:2070:G:C2	26:1H:2442:C:C2	3.04	0.46
40:B8:99:LEU:HB3	40:B8:101:PHE:HE1	1.77	0.46
26:1H:2632:A:O2'	26:1H:2811:G:O2'	2.27	0.46
1:1G:1053:G:O2'	1:1G:1054:C:P	2.74	0.46
1:1G:971:G:C6	1:1G:1364:U:O2'	2.69	0.46
26:1H:581:C:O2'	26:1H:582:G:H5'	2.15	0.46
26:14:2628:C:H1'	26:14:2781:A:H2'	1.97	0.46
9:8E:70:LYS:O	9:8E:74:ILE:HG13	2.16	0.46
46:D5:96:VAL:N	46:D5:128:VAL:O	2.41	0.46
26:1H:309:G:H4'	45:G8:18:GLY:HA2	1.98	0.46
2:12:141:GLU:O	2:12:145:LEU:HB2	2.15	0.46
26:1H:1843:C:H6	26:1H:1843:C:O5'	1.99	0.46
26:14:2080:G:O2'	26:14:2081:C:H5'	2.15	0.46
26:1H:2726:U:O2'	26:1H:2727:G:H8	1.99	0.46
30:39:152:GLU:HA	30:39:190:GLU:OE2	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
11:2I:44:SER:OG	11:2I:47:VAL:HG23	2.15	0.46
26:14:1665:A:H2'	26:14:1666:G:O4'	2.16	0.46
35:68:22:ILE:HD13	35:68:22:ILE:HA	1.59	0.46
30:31:117:ARG:HG2	30:31:192:LEU:HB2	1.98	0.46
28:11:218:ARG:HB3	28:11:219:PRO:HD2	1.98	0.46
1:1G:639:G:H2'	1:1G:640:A:H8	1.80	0.46
26:14:2212:A:O2'	26:14:2213:U:O5'	2.33	0.46
42:D8:21:ARG:HG2	42:D8:91:TYR:CE2	2.51	0.46
26:1H:2793:G:H2'	26:1H:2794:C:C6	2.51	0.46
50:H5:28:LEU:HA	50:H5:28:LEU:HD23	1.79	0.46
24:1L:9:A:H5'	24:1L:9:A:N3	2.31	0.46
45:C5:2:ARG:HD3	45:C5:2:ARG:HA	1.66	0.46
26:1H:2082:A:H2'	26:1H:2083:G:O4'	2.16	0.46
26:1H:1213:A:H1'	26:1H:1238:G:N3	2.30	0.46
1:13:32:A:H2'	1:13:33:A:C8	2.51	0.46
33:69:97:ILE:O	33:69:100:ALA:HB3	2.16	0.46
49:K8:64:LEU:HD22	49:K8:68:ARG:HD2	1.96	0.46
2:12:8:LYS:HE3	2:12:11:LEU:HD23	1.98	0.46
2:12:212:GLN:O	2:12:216:SER:N	2.34	0.46
1:13:1346:A:O3'	1:13:1347:G:H4'	2.16	0.46
26:1H:1359:A:N1	26:1H:1372:U:C4	2.83	0.46
26:14:2271:G:OP1	47:E5:18:ALA:HB1	2.15	0.46
24:3K:22:G:N7	24:3K:46:G:N2	2.63	0.46
45:C5:15:VAL:HG12	45:C5:21:LYS:HA	1.96	0.46
31:49:95:ARG:CG	31:49:96:ARG:H	2.27	0.46
55:M5:60:LEU:HB2	55:M5:61:LEU:H	1.38	0.46
19:AI:40:ILE:HD11	19:AI:62:ILE:HG23	1.97	0.46
1:1G:672:U:H2'	1:1G:673:G:H8	1.81	0.46
26:14:2420:C:OP1	55:M5:34:TRP:HB3	2.16	0.46
26:1H:1163:G:N3	26:1H:1164:G:C8	2.84	0.46
1:13:872:A:C4	1:13:874:G:N7	2.84	0.46
13:4I:3:ARG:HB3	13:4I:9:ILE:HG12	1.98	0.46
28:11:146:GLU:HB2	28:11:189:CYS:CB	2.44	0.46
2:1E:178:ARG:HD3	2:1E:178:ARG:HA	1.86	0.46
1:13:265:G:N2	1:13:267:C:H5'	2.31	0.46
1:1G:553:A:H2'	1:1G:554:C:H6	1.80	0.46
26:1H:1405:U:H2'	26:1H:1406:U:H6	1.80	0.46
39:A8:24:LEU:HB2	39:A8:85:VAL:CG1	2.46	0.46
29:29:199:ARG:HH12	29:29:202:LYS:NZ	2.14	0.46
26:1H:2261:C:O2'	26:1H:2262:U:H5'	2.15	0.46
3:2E:6:HIS:CD2	14:5I:49:HIS:HB3	2.50	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
37:88:12:GLN:HE21	37:88:72:LYS:HG3	1.81	0.46
1:1G:517:G:N3	1:1G:531:U:H5'	2.31	0.46
26:1H:1055:G:H1	26:1H:1104:C:H42	1.64	0.46
1:1G:1327:C:H2'	1:1G:1328:C:C6	2.51	0.46
19:AA:80:TYR:CZ	19:AA:82:GLY:HA2	2.50	0.46
26:14:579:G:H2'	26:14:580:C:H6	1.78	0.46
26:1H:1591:G:H2'	26:1H:1592:C:C6	2.51	0.46
32:51:87:LEU:HA	32:51:163:TYR:O	2.16	0.46
26:1H:85:G:OP2	45:G8:9:LYS:HB2	2.14	0.46
27:16:78:A:H2'	27:16:79:C:O4'	2.16	0.46
26:1H:6:A:H2'	26:1H:7:G:O4'	2.16	0.46
26:1H:950:G:C6	26:1H:951:C:C4	3.04	0.46
51:I5:57:GLU:O	51:I5:61:ARG:NH2	2.48	0.46
26:1H:2615:U:P	58:1H:3608:HOH:O	2.73	0.46
44:F8:50:LYS:O	44:F8:83:VAL:HA	2.16	0.46
1:1G:308:C:H2'	1:1G:309:G:H8	1.81	0.46
1:13:656:C:H2'	15:6I:28:GLN:NE2	2.31	0.46
26:14:1475:G:C2	26:14:1519:G:C2	3.04	0.46
28:19:236:GLY:N	28:19:237:GLU:HG2	2.31	0.46
15:6A:36:ILE:HD12	15:6A:63:ARG:HD3	1.98	0.46
26:14:1252:G:N3	41:85:33:ARG:HD2	2.30	0.46
26:14:560:C:O2	41:85:49:HIS:CE1	2.69	0.46
1:13:138:G:H1	1:13:225:C:H42	1.64	0.46
35:68:107:ARG:NH1	40:B8:36:GLU:HG2	2.31	0.46
49:G5:12:GLU:O	49:G5:16:LEU:HB3	2.15	0.46
1:1G:1134:G:C2	1:1G:1135:U:H1'	2.51	0.46
1:13:616:G:C2	1:13:617:G:N7	2.84	0.46
26:14:686:G:N7	54:L5:5:TRP:CH2	2.84	0.46
1:1G:1346:A:H3'	1:1G:1346:A:OP2	2.15	0.46
52:N8:9:LYS:HA	52:N8:9:LYS:HD3	1.74	0.46
2:12:10:LEU:HD13	2:12:10:LEU:HA	1.79	0.46
36:35:105:LEU:H	36:35:105:LEU:HD12	1.81	0.46
1:1G:1508:G:H1	1:1G:1527:C:H42	1.64	0.46
26:1H:2543:G:H1'	26:1H:2766:G:H5'	1.97	0.46
51:M8:16:CYS:SG	51:M8:36:CYS:N	2.89	0.46
2:1E:67:THR:N	2:1E:160:ASP:OD2	2.44	0.46
16:7I:71:ARG:O	16:7I:75:ARG:N	2.49	0.46
36:78:58:THR:O	36:78:62:LEU:HG	2.16	0.46
26:14:2275:C:H5'	26:14:2275:C:C6	2.49	0.46
30:31:63:LYS:HE2	30:31:67:GLN:HB3	1.98	0.46
30:31:64:ILE:HG23	30:31:65:TRP:NE1	2.31	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:1H:2344:U:O2'	53:O8:37:ARG:HG2	2.15	0.46
42:95:38:LEU:HD22	42:95:55:ALA:HB1	1.99	0.46
29:21:18:ASP:CB	40:B8:82:LEU:HD11	2.46	0.46
1:13:991:U:O2'	1:13:992:U:O5'	2.33	0.46
26:14:1426:G:H5''	26:14:1427:A:OP2	2.16	0.46
26:14:2299:G:N1	26:14:2318:G:C8	2.84	0.46
1:13:575:G:C5	1:13:881:G:C2	3.04	0.46
37:45:85:LYS:HG2	37:45:86:GLY:H	1.81	0.46
26:14:2115:G:N2	26:14:2172:U:H3	2.14	0.46
12:3I:84:LEU:HB2	12:3I:105:TYR:HE2	1.81	0.46
36:78:112:LEU:O	36:78:128:HIS:HB2	2.16	0.46
24:3L:58:A:H1'	24:3L:60:U:C5	2.51	0.46
26:1H:773:U:H5'	28:11:47:GLY:HA3	1.98	0.46
37:45:59:ARG:H	37:45:59:ARG:HG3	1.44	0.46
1:1G:1163:C:N3	1:1G:1174:G:N2	2.64	0.46
42:95:29:PRO:HA	42:95:61:VAL:CG1	2.46	0.46
7:6E:62:PHE:HA	7:6E:124:LEU:HD21	1.97	0.46
13:4I:15:VAL:O	13:4I:19:LEU:HD23	2.16	0.46
2:12:189:ASP:H	2:12:192:SER:HB2	1.81	0.46
46:D5:92:SER:O	46:D5:94:GLU:N	2.48	0.46
26:1H:813:U:H2'	26:1H:814:C:C6	2.51	0.46
26:1H:2016:U:H1'	52:N8:6:VAL:HG13	1.98	0.46
19:AA:35:SER:O	19:AA:71:LEU:HD12	2.16	0.46
26:14:2340:G:H2'	26:14:2341:G:H8	1.80	0.46
10:1I:3:LYS:N	10:1I:75:ILE:HA	2.31	0.46
26:14:1421:G:C2	26:14:1422:G:N7	2.83	0.46
6:52:76:ALA:O	6:52:80:ARG:HG3	2.16	0.46
4:32:28:SER:HA	4:32:29:PRO:HA	1.61	0.46
14:5A:26:ARG:CZ	14:5A:47:LEU:HD21	2.46	0.46
36:35:3:LEU:HD12	36:35:3:LEU:H	1.80	0.45
26:14:1113:U:OP1	32:59:3:ARG:N	2.49	0.45
26:1H:1515:C:H2'	26:1H:1516:U:H6	1.81	0.45
1:1G:1298:C:H5	7:62:114:ARG:HD2	1.81	0.45
46:D5:157:LEU:HA	46:D5:158:PRO:HD2	1.53	0.45
2:12:102:LEU:HD21	2:12:162:ILE:HG21	1.98	0.45
1:1G:1320:C:H2'	1:1G:1321:C:C6	2.51	0.45
10:1A:49:VAL:O	10:1A:60:ARG:HB2	2.15	0.45
26:14:690:G:H2'	26:14:691:C:C6	2.51	0.45
1:1G:29:G:H5'	1:1G:296:U:OP1	2.16	0.45
26:1H:570:G:H5''	58:1H:3815:HOH:O	2.16	0.45
1:1G:406:G:H1'	1:1G:495:A:C6	2.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1G:279:A:C8	1:1G:281:G:C2	3.04	0.45
35:68:120:GLU:HB2	40:B8:68:TYR:HE2	1.81	0.45
3:2E:3:ASN:O	3:2E:4:LYS:HG2	2.16	0.45
26:14:548:A:C6	26:14:549:G:H1'	2.51	0.45
26:1H:2294:C:O2'	26:1H:2295:C:H5'	2.16	0.45
26:1H:208:C:H2'	26:1H:209:C:H6	1.81	0.45
46:H8:129:SER:HA	46:H8:130:PRO:HD3	1.77	0.45
1:1G:1233:G:H2'	1:1G:1234:C:C6	2.51	0.45
2:1E:58:ILE:HB	2:1E:221:LEU:HD12	1.97	0.45
26:14:2648:C:H2'	26:14:2649:U:C6	2.51	0.45
1:13:665:A:N3	1:13:732:C:H2'	2.31	0.45
26:14:696:G:H2'	26:14:697:C:H6	1.80	0.45
1:1G:954:G:H2'	1:1G:955:U:C6	2.51	0.45
35:68:64:ARG:HG2	35:68:79:PHE:CG	2.50	0.45
23:2L:11:A:H8	23:2L:11:A:O5'	1.99	0.45
7:62:63:LYS:HG3	7:62:64:GLN:N	2.31	0.45
30:39:178:PRO:HB3	30:39:198:ALA:HB1	1.98	0.45
26:14:270(G):C:H2'	26:14:270(H):C:C6	2.51	0.45
39:65:41:ASP:CG	39:65:44:LYS:HD2	2.36	0.45
36:78:17:LYS:HE2	36:78:27:HIS:CD2	2.51	0.45
26:1H:1899:G:N2	26:1H:1902:C:H5	2.13	0.45
26:14:662:G:H5'	36:35:15:ARG:CA	2.34	0.45
26:1H:1728:G:H3'	26:1H:1729:A:C5'	2.46	0.45
26:1H:1478:G:H1'	26:1H:1557:C:O2'	2.17	0.45
1:13:1022:G:H2'	1:13:1023:G:H8	1.82	0.45
4:32:121:VAL:O	4:32:134:ASP:HA	2.16	0.45
45:C5:82:PRO:CB	45:C5:97:ARG:HB3	2.46	0.45
43:E8:70:TYR:CD1	43:E8:70:TYR:N	2.83	0.45
1:1G:457:C:H2'	1:1G:458:C:C6	2.51	0.45
30:31:64:ILE:HA	30:31:64:ILE:HD13	1.68	0.45
4:3E:9:CYS:HB3	4:3E:32:ALA:HB2	1.98	0.45
1:1G:1335:C:OP1	1:1G:1337:G:N2	2.35	0.45
26:1H:2287:A:C2	26:1H:2346:A:H2	2.34	0.45
1:1G:1166:G:N2	1:1G:1170:A:OP2	2.49	0.45
47:I8:11:ARG:HG3	47:I8:11:ARG:H	1.44	0.45
1:13:1013:G:N2	1:13:1016:A:OP2	2.47	0.45
44:F8:57:LEU:HG	44:F8:78:LYS:HG2	1.99	0.45
1:1G:75:C:H2'	1:1G:76:G:O4'	2.16	0.45
3:2E:150:LYS:HG3	3:2E:169:ALA:HB2	1.98	0.45
36:35:99:LEU:HD12	36:35:102:ARG:NH2	2.31	0.45
48:F5:83:GLU:N	48:F5:83:GLU:OE1	2.49	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:2467:C:H4'	37:45:123:HIS:CD2	2.52	0.45
36:78:114:ILE:O	36:78:114:ILE:HG13	2.16	0.45
36:78:83:VAL:O	36:78:114:ILE:HA	2.16	0.45
29:29:42:ASP:HB3	29:29:44:TYR:CE2	2.51	0.45
38:55:56:LYS:HD2	38:55:88:ARG:HA	1.98	0.45
1:1G:748:C:O5'	1:1G:748:C:H6	1.99	0.45
1:1G:1443:G:N2	40:75:119:LYS:HB2	2.31	0.45
34:58:41:ASP:N	34:58:41:ASP:OD1	2.50	0.45
13:4I:15:VAL:HG23	13:4I:43:THR:O	2.16	0.45
13:4I:19:LEU:O	13:4I:22:ILE:HG13	2.17	0.45
17:8A:100:LYS:HD3	17:8A:100:LYS:HA	1.51	0.45
26:1H:902:C:H2'	26:1H:903:C:H6	1.81	0.45
30:31:7:TYR:O	30:31:22:ALA:N	2.49	0.45
2:1E:76:GLN:NE2	2:1E:207:ALA:H	2.13	0.45
38:98:53:HIS:HB2	38:98:94:TYR:HE2	1.81	0.45
32:59:97:ARG:HG2	32:59:98:LEU:H	1.81	0.45
54:L5:34:ARG:NH1	54:L5:39:ARG:HG3	2.32	0.45
30:31:133:ASN:HB3	30:31:138:GLU:OE1	2.15	0.45
4:32:155:LEU:O	4:32:158:ILE:HG13	2.16	0.45
4:32:97:LEU:O	4:32:100:ARG:HB2	2.16	0.45
1:1G:1151:A:O2'	1:1G:1152:A:O5'	2.34	0.45
32:59:136:ILE:H	32:59:136:ILE:HG13	1.48	0.45
33:61:79:ILE:HD13	33:61:79:ILE:HA	1.76	0.45
2:12:172:ILE:H	2:12:172:ILE:HD12	1.81	0.45
1:1G:979:C:OP1	1:1G:1223:C:N4	2.49	0.45
26:14:176:G:O2'	26:14:177:G:H5'	2.16	0.45
26:1H:2505:G:H2'	26:1H:2576:G:O6	2.15	0.45
1:1G:1145:C:H1'	1:1G:1146:A:N7	2.32	0.45
41:85:92:ARG:O	41:85:94:ASN:N	2.48	0.45
30:39:25:PRO:HB2	30:39:27:GLU:HB2	1.98	0.45
26:14:2134:A:H2	26:14:2159:G:HO2'	1.63	0.45
21:1B:10:ARG:HA	21:1B:13:ILE:HD12	1.97	0.45
1:13:1129:C:N4	1:13:1139:G:H1	2.13	0.45
26:14:2261:C:H1'	26:14:2388:A:N3	2.31	0.45
26:1H:1050:A:H2'	26:1H:1051:G:O4'	2.17	0.45
3:2E:16:ARG:HA	3:2E:16:ARG:HD2	1.69	0.45
1:1G:1028(A):C:N4	1:1G:1032(B):G:H22	2.14	0.45
26:14:2296:U:H4'	26:14:2297:C:OP1	2.16	0.45
1:13:955:U:H1'	1:13:1227:A:N6	2.30	0.45
26:1H:2784:C:H1'	29:21:37:ARG:NH1	2.30	0.45
9:8E:27:THR:HB	9:8E:62:TYR:HA	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:1E:12:GLU:HA	2:1E:16:HIS:CG	2.52	0.45
26:14:528:A:OP2	34:15:114:ARG:NH1	2.49	0.45
26:1H:1419:A:C8	26:1H:1421:G:C6	3.05	0.45
26:14:2346:A:H5''	26:14:2383:G:H1'	1.97	0.45
49:K8:23:LYS:HE3	49:K8:27:GLU:OE2	2.16	0.45
53:O8:17:LYS:O	53:O8:18:ARG:HB2	2.16	0.45
26:1H:535:C:O3'	41:C8:53:ARG:NH1	2.48	0.45
39:65:86:ALA:O	39:65:87:PHE:HB2	2.17	0.45
26:14:863:A:H2'	26:14:864:G:H8	1.82	0.45
26:14:1257:C:H4'	30:39:83:PHE:CE1	2.52	0.45
37:88:11:LYS:CE	37:88:86:GLY:HA2	2.46	0.45
26:1H:1155:A:H5'	41:C8:55:ARG:NE	2.30	0.45
38:98:2:ARG:HB3	38:98:3:HIS:H	1.48	0.45
1:1G:35:G:C2	1:1G:550:G:C2	3.03	0.45
26:14:354:G:H2'	26:14:355:G:C8	2.52	0.45
34:15:29:LYS:H	34:15:29:LYS:HG2	1.38	0.45
32:59:19:VAL:HG12	32:59:20:ALA:H	1.81	0.45
26:14:931:G:H3'	26:14:931:G:H8	1.81	0.45
1:13:684:A:H2'	1:13:685:G:C8	2.51	0.45
20:BI:33:ILE:O	20:BI:37:SER:OG	2.33	0.45
39:65:99:LYS:NZ	39:65:103:GLU:OE1	2.36	0.45
8:7E:25:ASP:OD1	8:7E:60:ARG:HG3	2.16	0.45
43:E8:40:ASN:O	43:E8:41:LYS:HG2	2.16	0.45
26:14:2716:U:O2'	26:14:2717:G:H5'	2.16	0.45
26:1H:2795:G:H3'	26:1H:2797:U:C5'	2.46	0.45
31:49:144:ILE:HG22	31:49:146:TYR:H	1.82	0.45
17:8A:40:LYS:HD3	17:8A:42:TYR:CE1	2.51	0.45
28:19:94:LEU:HD23	28:19:94:LEU:HA	1.73	0.45
14:5A:53:LEU:HD23	14:5A:53:LEU:HA	1.67	0.45
19:AA:22:LEU:O	19:AA:27:GLU:HA	2.15	0.45
49:G5:63:VAL:O	49:G5:67:LYS:HG2	2.15	0.45
26:14:270(L):U:O2'	26:14:270(M):U:OP1	2.32	0.45
9:8E:112:LYS:HD3	9:8E:113:LYS:N	2.31	0.45
26:1H:1385:G:O6	26:1H:1403:C:N4	2.50	0.45
55:Q8:45:GLY:CA	55:Q8:46:ARG:C	2.84	0.45
26:1H:1416:G:H2'	26:1H:1417:C:C5	2.51	0.45
26:1H:2159:G:H2'	26:1H:2160:G:O4'	2.17	0.45
27:1J:65:C:H41	27:1J:108:C:C2'	2.29	0.45
26:1H:654(C):G:H2'	26:1H:654(D):G:O4'	2.16	0.45
20:BI:69:GLY:O	20:BI:73:HIS:CE1	2.70	0.45
31:49:66:GLN:NE2	31:49:93:THR:O	2.49	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:2496:C:P	37:45:81:VAL:HG12	2.56	0.45
37:88:119:ARG:HB3	37:88:119:ARG:HE	1.42	0.45
1:1G:1057:G:H2'	1:1G:1058:G:C8	2.52	0.45
26:14:2002:G:C5	58:14:3900:HOH:O	2.68	0.45
44:F8:12:VAL:HG13	44:F8:27:THR:O	2.16	0.45
26:14:1342:A:H2	26:14:1602:U:N3	2.12	0.45
4:3E:188:LEU:HA	4:3E:189:PRO:HD3	1.69	0.45
1:1G:421:U:O2'	1:1G:423:G:N7	2.49	0.45
7:62:93:PRO:HG2	7:62:94:ARG:HD3	1.98	0.45
1:1G:980:C:H5'	1:1G:981:U:OP2	2.17	0.45
26:14:1794:U:H2'	26:14:1795:C:C6	2.51	0.45
28:19:16:MET:HG3	28:19:206:LEU:O	2.16	0.45
28:19:206:LEU:HD22	28:19:211:ARG:HG2	1.98	0.45
47:I8:42:GLY:C	47:I8:57:PHE:HD2	2.20	0.45
1:1G:793:U:O4	1:1G:1517:G:H5''	2.17	0.45
1:13:1315:U:C5	1:13:1316:G:C5	3.05	0.45
26:14:1171:G:N2	26:14:1178:C:H42	2.14	0.45
7:6E:50:ILE:O	7:6E:54:THR:HG23	2.17	0.45
1:1G:16:A:H2'	1:1G:17:U:C6	2.51	0.45
1:1G:991:U:O2	1:1G:993:G:H8	2.00	0.45
45:C5:88:LYS:O	45:C5:89:PHE:HB3	2.15	0.45
1:13:465:A:N7	1:13:467:G:C6	2.84	0.45
47:I8:50:ASN:HD22	47:I8:83:PRO:HD3	1.80	0.45
26:14:1449:A:H5'	26:14:1449(A):G:OP2	2.16	0.45
46:D5:175:VAL:HA	46:D5:177:PRO:HD3	1.97	0.45
1:1G:187:C:H2'	1:1G:188:U:O4'	2.16	0.45
26:1H:784:A:C5	28:11:229:VAL:HG21	2.51	0.45
1:1G:49:U:C2	1:1G:361:G:N2	2.85	0.45
26:14:2511:U:H2'	26:14:2512:C:C6	2.51	0.45
8:7E:134:ILE:HG22	8:7E:135:CYS:SG	2.57	0.45
40:75:85:LYS:HD2	40:75:87:ASP:OD1	2.17	0.45
29:29:15:PHE:CE1	29:29:20:ALA:HB2	2.52	0.45
7:6E:26:PHE:CE2	7:6E:30:ILE:HD11	2.52	0.45
26:14:2754:U:H5''	26:14:2754:U:H6	1.82	0.45
26:14:1572:A:H8	26:14:1572:A:O5'	1.99	0.45
34:15:96:GLU:H	34:15:96:GLU:CD	2.18	0.45
26:14:1359:A:N7	26:14:1372:U:O4	2.49	0.45
26:1H:1665:A:H2'	26:1H:1666:G:O4'	2.17	0.45
1:13:112:G:P	16:7I:27:LYS:HD2	2.57	0.45
4:3E:59:ARG:HH21	4:3E:66:ARG:HH12	1.64	0.45
1:13:1183:A:O2'	1:13:1184:G:OP1	2.28	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:32:22:LYS:HG3	4:32:26:CYS:SG	2.56	0.45
27:1J:40:U:C2'	27:1J:45:A:H61	2.30	0.45
26:1H:2789:C:H1'	26:1H:2892:A:C2	2.50	0.45
46:H8:76:LEU:HA	46:H8:83:PRO:HA	1.98	0.45
21:1B:9:ARG:HG3	21:1B:10:ARG:HG3	1.97	0.45
13:4I:3:ARG:CZ	13:4I:7:VAL:HG13	2.47	0.45
1:13:1309:G:C6	1:13:1329:A:C2	3.04	0.45
55:M5:54:GLU:HA	55:M5:57:ARG:NH1	2.32	0.45
1:1G:564:C:HO2'	8:72:91:ARG:HH22	1.61	0.45
36:35:125:VAL:HG13	36:35:144:GLU:HB3	1.99	0.45
34:15:134:ARG:HB3	34:15:134:ARG:HE	1.49	0.45
26:1H:1069:A:H4'	26:1H:1070:A:H5''	1.99	0.45
1:1G:1169:A:H2'	1:1G:1170:A:C8	2.52	0.45
26:14:1533:C:H42	26:14:1538:G:H1	1.64	0.45
23:2K:63:C:O2	23:2K:64:G:C8	2.70	0.45
26:1H:1486:A:C2	26:1H:1487:G:C5	3.04	0.45
26:1H:1348:G:C2'	26:1H:1349:A:H5''	2.46	0.45
26:1H:931:G:H4'	50:L8:24:LYS:HZ3	1.81	0.45
43:E8:57:ASN:O	43:E8:62:HIS:HD2	2.00	0.45
30:39:181:LEU:CD2	30:39:186:ILE:HD11	2.46	0.45
26:1H:479:A:N3	26:1H:481:G:H5'	2.32	0.45
24:1L:14:A:N1	24:1L:22:G:H1'	2.31	0.45
10:1I:78:ASN:O	10:1I:81:THR:N	2.50	0.45
46:D5:52:SER:O	46:D5:54:HIS:N	2.50	0.45
1:1G:678:U:H2'	1:1G:679:C:C6	2.51	0.45
1:13:417:C:H2'	1:13:418:C:H6	1.80	0.45
3:2E:121:ALA:O	3:2E:125:GLU:HG3	2.17	0.45
26:14:225:A:N6	26:14:226:G:C2	2.85	0.45
16:7A:4:ILE:HB	16:7A:66:PRO:HB3	1.98	0.45
1:1G:1420:C:H6	1:1G:1420:C:O5'	2.00	0.45
3:22:175:LEU:H	3:22:175:LEU:HD12	1.82	0.45
1:13:967:C:O5'	1:13:967:C:H6	2.00	0.45
35:68:53:LYS:HA	35:68:53:LYS:HD2	1.82	0.45
36:35:41:ARG:HD2	36:35:41:ARG:N	2.31	0.45
26:14:1357:U:H2'	26:14:1358:G:O4'	2.17	0.45
31:41:112:PRO:HB3	51:M8:37:SER:N	2.12	0.45
1:1G:1149:C:H2'	1:1G:1150:U:O4'	2.17	0.45
1:13:922:G:C6	1:13:923:A:C6	3.04	0.45
26:14:1050:A:N3	26:14:2751:G:H2'	2.32	0.45
55:Q8:45:GLY:HA2	55:Q8:46:ARG:C	2.36	0.45
26:1H:2502:G:N7	58:1H:3791:HOH:O	2.36	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:12:75:LYS:O	2:12:75:LYS:HD2	2.16	0.45
45:C5:17:SER:CB	45:C5:71:LYS:HB3	2.44	0.45
26:1H:654(H):G:H2'	26:1H:654(H):G:N3	2.32	0.45
26:1H:1534:G:N2	26:1H:1538:G:H1	2.14	0.45
26:1H:1465:G:H2'	26:1H:1466:G:H8	1.81	0.45
1:13:633:G:OP2	1:13:633:G:H8	2.00	0.45
26:14:1826:G:H4'	28:19:242:ARG:HE	1.80	0.45
23:2K:48:U:OP2	23:2K:48:U:C6	2.69	0.45
43:A5:73:ALA:HB3	43:A5:106:ILE:CD1	2.43	0.45
44:F8:12:VAL:HG22	44:F8:17:ALA:HB2	1.99	0.45
2:1E:70:PHE:HD2	2:1E:91:PRO:O	2.00	0.45
26:14:1011:G:N3	26:14:1151:G:C2	2.84	0.45
38:98:42:LYS:O	38:98:45:ARG:HD2	2.17	0.45
23:2L:32:G:C4	23:2L:33:OMC:C5	3.04	0.45
47:I8:63:VAL:HG23	47:I8:64:ASP:O	2.15	0.45
1:13:191(C):G:H2'	1:13:191(D):U:C6	2.52	0.45
26:1H:1444:G:N2	26:1H:1548:C:N3	2.65	0.45
31:49:110:ALA:HA	31:49:140:ILE:O	2.16	0.45
26:1H:1280:G:N2	26:1H:1291:C:C2	2.85	0.45
45:C5:87:LYS:HG2	45:C5:88:LYS:N	2.31	0.45
26:14:1542:G:H3'	26:14:1543:A:H5''	1.99	0.45
18:9I:41:LYS:HE3	18:9I:41:LYS:HB3	1.73	0.45
51:M8:15:ILE:HG22	51:M8:20:ASN:OD1	2.16	0.45
38:98:29:LEU:HD23	38:98:70:LEU:HD11	1.99	0.45
26:14:768:G:H2'	26:14:769:G:H8	1.81	0.45
2:1E:130:ARG:HB2	2:1E:134:GLU:HB2	1.98	0.45
1:13:1239:A:H62	1:13:1299:A:H62	1.64	0.45
28:11:121:PRO:HB3	28:11:135:PHE:CE2	2.52	0.45
27:1J:111:U:H2'	27:1J:112:G:H8	1.82	0.45
1:1G:1104:G:H4'	2:12:111:ARG:HE	1.82	0.45
26:14:550:G:O2'	26:14:1220:A:N3	2.44	0.45
53:K5:51:GLU:HG2	53:K5:52:VAL:N	2.30	0.45
41:85:110:VAL:O	41:85:114:LYS:HG2	2.16	0.45
46:D5:19:ARG:NH1	46:D5:84:GLU:HB2	2.32	0.45
32:51:95:ARG:NH1	32:51:95:ARG:HB3	2.31	0.45
44:B5:60:ARG:HG2	44:B5:60:ARG:HH11	1.81	0.45
31:41:43:LEU:HD12	31:41:43:LEU:HA	1.61	0.45
33:69:113:ARG:HA	33:69:113:ARG:HD3	1.73	0.45
23:2L:36:A:C2	25:4L:18:G:C2	3.05	0.45
10:1A:82:ILE:O	10:1A:86:MET:HB2	2.16	0.45
28:19:75:ILE:O	28:19:118:VAL:HG23	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:78:19:VAL:HG12	36:78:21:ARG:N	2.11	0.45
1:13:1366:C:H2'	1:13:1367:C:H6	1.81	0.45
26:14:1204:A:C2	26:14:1241:A:N1	2.85	0.45
26:14:1042:G:H2'	26:14:1043:C:C6	2.52	0.45
26:14:1034:G:H2'	26:14:1035:U:O4'	2.17	0.45
26:1H:2592:G:C6	26:1H:2593:U:N3	2.84	0.45
1:1G:854:G:N1	1:1G:855:G:N7	2.64	0.45
33:69:73:GLU:HG2	33:69:137:PRO:O	2.17	0.45
26:14:2712(A):A:H5''	26:14:2713:A:OP2	2.17	0.45
51:M8:13:ARG:HA	51:M8:24:THR:HG21	1.99	0.45
38:98:83:ILE:HG22	38:98:87:TYR:CE2	2.46	0.45
26:14:270(S):G:OP1	48:F5:76:ARG:NH1	2.49	0.45
1:1G:1239:A:H4'	1:1G:1240:U:H5'	1.99	0.45
26:1H:1446:C:H2'	26:1H:1447:G:H8	1.81	0.45
1:1G:538:G:H2'	1:1G:539:A:H8	1.81	0.45
26:1H:583:G:H5''	41:C8:10:ARG:NH1	2.31	0.45
37:88:72:LYS:HB3	37:88:94:VAL:HG23	1.97	0.45
30:39:40:GLN:OE1	30:39:182:ASN:HB2	2.17	0.45
1:13:674:G:H2'	1:13:675:A:H8	1.81	0.45
30:39:107:LYS:HD2	30:39:205:ARG:HG3	1.98	0.45
32:51:92:ILE:HD12	32:51:93:GLY:H	1.81	0.45
26:14:1153:C:OP1	41:85:93:LYS:NZ	2.50	0.45
24:1L:57:G:H2'	24:1L:58:A:H5''	1.98	0.45
26:14:108:U:H2'	26:14:109:G:C8	2.52	0.45
26:1H:26:G:C6	26:1H:27:G:N1	2.85	0.45
1:13:1207:G:H2'	1:13:1208:C:C6	2.52	0.45
1:13:358:U:H2'	1:13:359:U:O4'	2.17	0.45
26:14:2059:A:H5''	26:14:2060:A:OP2	2.17	0.45
39:A8:72:ALA:O	39:A8:76:LYS:HG3	2.16	0.45
4:32:74:GLN:O	4:32:78:LEU:HD12	2.17	0.45
26:1H:1392:A:C6	26:1H:1393:A:N1	2.84	0.45
35:68:60:ALA:HB1	35:68:84:ALA:HB1	1.99	0.45
15:6I:82:ILE:HG22	15:6I:83:GLU:N	2.32	0.45
26:1H:183:C:H42	26:1H:213:A:H61	1.64	0.45
21:1F:6:ARG:HG3	21:1F:6:ARG:H	1.41	0.45
1:13:730:G:C5	1:13:731:G:H1'	2.52	0.45
4:3E:59:ARG:NH2	4:3E:66:ARG:HH12	2.15	0.45
1:1G:1063:C:H3'	1:1G:1064:G:H2'	1.99	0.45
26:14:921:G:H2'	26:14:922:U:C6	2.52	0.45
49:K8:15:LYS:HE2	49:K8:67:LYS:HZ3	1.81	0.45
26:14:2332:U:H5'	47:E5:43:THR:HG21	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1G:1345:U:O2	1:1G:1377:A:N6	2.50	0.45
1:1G:942:G:C2	1:1G:1342:C:C2	3.05	0.45
9:82:125:TYR:HD1	9:82:126:SER:N	2.15	0.45
40:75:50:ILE:HD11	40:75:102:ILE:CD1	2.45	0.45
1:1G:498:A:H4'	1:1G:500:G:OP1	2.16	0.45
29:21:171:GLU:OE1	29:21:185:LYS:HD3	2.17	0.45
29:21:34:VAL:HG21	29:21:77:ILE:HD13	1.99	0.45
26:14:1412:A:H2'	26:14:1413:G:H8	1.81	0.45
1:1G:567:G:O6	12:3A:5:PRO:HD3	2.17	0.45
1:13:1015:A:H1'	1:13:1219:U:H5'	1.98	0.45
1:13:1248:A:C2	9:8E:70:LYS:HD2	2.52	0.45
20:BI:100:ILE:HG13	20:BI:101:GLY:H	1.82	0.45
26:14:1025:G:O2'	26:14:1026:U:OP1	2.27	0.45
45:C5:52:SER:H	45:C5:57:GLN:N	2.14	0.45
32:51:86:GLU:HG2	32:51:87:LEU:H	1.82	0.45
26:1H:806:C:H2'	26:1H:807:U:H6	1.82	0.45
55:Q8:38:GLY:HA2	55:Q8:39:LYS:C	2.38	0.45
26:1H:2068:U:N3	26:1H:2430:A:C2	2.74	0.45
26:1H:302:C:H2'	26:1H:303:U:C6	2.52	0.45
29:21:92:THR:O	29:21:95:ILE:HG22	2.17	0.45
8:7E:82:HIS:HB3	8:7E:138:TRP:CH2	2.52	0.45
26:1H:775:G:C4	26:1H:794:G:C8	3.05	0.45
16:7A:9:PHE:HB2	16:7A:16:HIS:O	2.17	0.45
26:1H:556:G:H2'	26:1H:557:U:H6	1.82	0.45
4:32:126:ILE:HG22	4:32:127:THR:N	2.32	0.45
30:39:165:ARG:HH11	30:39:165:ARG:HB3	1.81	0.45
30:31:135:LYS:HB3	30:31:138:GLU:HG3	1.98	0.45
32:51:95:ARG:HH11	32:51:95:ARG:HB3	1.82	0.45
26:1H:992:C:OP1	42:D8:74:LYS:NZ	2.50	0.45
43:A5:18:ARG:HG3	43:A5:76:VAL:HG13	1.99	0.45
1:13:753:A:H4'	1:13:754:C:H5''	1.99	0.45
7:6E:80:VAL:C	7:6E:82:GLY:H	2.20	0.45
26:1H:2666:C:H5''	26:1H:2667:C:OP2	2.16	0.45
40:B8:51:ARG:HB2	40:B8:98:LYS:HD3	1.98	0.45
26:1H:289:A:H2'	26:1H:290:G:O4'	2.17	0.45
1:13:134:A:H1'	1:13:325:A:C5	2.52	0.45
47:E5:56:ASP:CG	47:E5:58:THR:HG1	2.20	0.45
48:F5:44:PRO:HB2	48:F5:46:LEU:HD12	1.98	0.45
43:E8:29:LEU:HD21	43:E8:33:ARG:NH2	2.32	0.45
23:2L:29:C:H6	23:2L:29:C:O5'	1.99	0.45
33:61:9:LEU:HA	33:61:9:LEU:HD12	1.69	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:32:146:ILE:H	4:32:146:ILE:HD12	1.81	0.45
1:13:181:G:HO2'	1:13:182:U:H6	1.61	0.45
1:1G:704:A:H5'	1:1G:705:U:OP2	2.17	0.45
26:1H:28:A:C2	26:1H:513:A:C8	3.04	0.45
43:E8:2:GLU:HB2	43:E8:107:LEU:O	2.17	0.45
12:3A:38:THR:OG1	12:3A:57:LYS:HB2	2.17	0.45
36:78:18:ARG:C	36:78:19:VAL:HG22	2.37	0.45
27:1J:15:A:H1'	27:1J:109:G:C4	2.52	0.45
1:1G:1145:C:H4'	1:1G:1146:A:O5'	2.17	0.45
9:82:17:VAL:HG21	9:82:80:GLY:C	2.37	0.45
1:13:926:G:H5'	1:13:927:G:O5'	2.16	0.45
1:1G:1505:G:H4'	1:1G:1506:U:H5''	1.99	0.45
1:1G:795:C:H1'	1:1G:1506:U:C5	2.52	0.45
39:A8:36:TYR:N	39:A8:36:TYR:HD1	2.13	0.45
26:14:1113:U:O5'	26:14:1113:U:H6	1.99	0.45
1:1G:1043:C:H2'	1:1G:1044:A:H8	1.82	0.45
26:1H:1163:G:C2	26:1H:1164:G:N7	2.85	0.45
40:B8:30:VAL:HG23	40:B8:83:ILE:HG23	1.98	0.45
30:31:66:PRO:O	30:31:67:GLN:CB	2.64	0.45
30:31:127:GLU:HA	30:31:127:GLU:OE2	2.11	0.45
13:4I:94:ARG:HD3	13:4I:94:ARG:HA	1.70	0.45
22:1K:27:G:H2'	22:1K:28:G:H8	1.82	0.45
26:1H:2058:A:N6	58:1H:3615:HOH:O	2.48	0.45
1:13:1103:C:H2'	1:13:1104:G:O4'	2.17	0.45
1:1G:1240:U:C2	7:62:32:ARG:HG3	2.51	0.45
29:29:197:ILE:O	29:29:197:ILE:HG13	2.17	0.45
5:4E:41:VAL:HG13	5:4E:113:ALA:HB2	1.98	0.45
26:1H:1858:G:OP2	26:1H:1858:G:H8	2.00	0.45
1:1G:1300:G:HO2'	1:1G:1301:U:P	2.38	0.45
26:14:1678:G:N2	26:14:1989:G:N2	2.65	0.45
26:1H:919:G:H4'	27:16:81:G:H4'	1.98	0.45
51:I5:18:CYS:N	51:I5:19:GLY:HA2	2.29	0.45
1:1G:1325:C:H2'	1:1G:1326:C:C6	2.52	0.45
35:68:71:ARG:NH2	35:68:122:LEU:O	2.48	0.45
22:1K:52:G:H2'	22:1K:53:G:O4'	2.17	0.45
26:1H:2040:C:H2'	26:1H:2041:U:O4'	2.17	0.45
26:14:840:C:H42	26:14:938:G:H1	1.65	0.45
26:1H:1637:A:H4'	26:1H:2711:A:O2'	2.17	0.45
5:4E:11:ILE:HG12	5:4E:31:LEU:HB3	1.98	0.45
1:13:456:C:H42	1:13:476:G:H1	1.65	0.45
36:78:39:LYS:HG3	36:78:45:LEU:CD2	2.47	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:1H:1125:G:OP2	26:1H:1126:A:O2'	2.32	0.45
32:51:92:ILE:HD11	32:51:160:LYS:NZ	2.32	0.45
17:8A:59:ILE:HG22	17:8A:71:PHE:HD2	1.81	0.45
31:49:114:ILE:HG12	31:49:140:ILE:HG21	1.99	0.45
26:1H:2093:G:C6	26:1H:2225:A:C8	3.05	0.45
1:1G:1152:A:H2'	1:1G:1153:C:H6	1.81	0.45
26:1H:2667:C:H1'	32:51:109:PHE:HD1	1.81	0.45
26:1H:1045:A:H1'	26:1H:1047:G:N3	2.32	0.45
48:J8:97:LEU:HG	48:J8:98:LEU:H	1.82	0.45
1:13:1044:A:C5	1:13:1045:C:H1'	2.52	0.45
38:55:10:LEU:HD23	38:55:10:LEU:HA	1.71	0.45
26:1H:2308:G:N3	26:1H:2308:G:H2'	2.32	0.45
26:1H:2592:G:C5	26:1H:2593:U:C4	3.05	0.45
22:1K:76:A:HO3'	26:1H:2506:U:H1'	1.82	0.45
22:1K:76:A:H2	26:1H:2602:A:C2	2.34	0.45
4:3E:62:GLN:O	4:3E:66:ARG:HD3	2.16	0.45
32:51:83:TYR:HA	32:51:135:GLY:H	1.82	0.45
1:1G:191(F):U:H2'	1:1G:191:G:C8	2.52	0.45
2:12:54:THR:HG23	2:12:199:TYR:HB3	1.99	0.45
30:39:122:LYS:HB3	30:39:191:ARG:HB2	1.98	0.45
34:58:22:THR:HG22	34:58:23:LEU:N	2.32	0.45
45:G8:97:ARG:HB3	45:G8:99:CYS:SG	2.57	0.45
26:1H:1799:G:O2'	26:1H:1800:C:OP2	2.24	0.45
1:1G:827:U:H3	1:1G:872:A:H62	1.63	0.45
23:2K:20:G:N2	23:2K:58:A:H1'	2.32	0.45
30:39:53:THR:HG22	30:39:56:GLU:CG	2.45	0.45
35:68:71:ARG:HH21	35:68:77:ILE:HG21	1.82	0.45
2:12:19:HIS:CE1	2:12:206:ASP:HB2	2.51	0.45
12:3I:82:VAL:HG13	12:3I:105:TYR:HB3	1.97	0.45
1:1G:300:A:H1'	1:1G:565:U:O2	2.17	0.45
41:85:66:ASN:CB	41:85:76:TYR:HB2	2.47	0.45
26:1H:2109:U:H2'	26:1H:2110:G:C8	2.52	0.45
1:1G:677:U:H3	1:1G:713:G:H22	1.64	0.45
16:7I:78:GLY:HA2	16:7I:81:ARG:HH12	1.82	0.45
1:13:689:C:C4	1:13:690:G:C2	3.05	0.45
45:G8:40:GLU:HB3	45:G8:64:GLU:OE1	2.17	0.45
26:14:2302:G:N2	26:14:2314:C:O2	2.48	0.45
3:22:18:TRP:HE1	14:5A:55:GLY:N	2.15	0.45
11:2I:125:PHE:O	11:2I:126:ARG:HG2	2.17	0.45
1:1G:857:C:H2'	1:1G:858:G:O4'	2.17	0.45
23:2L:54:G:H2'	23:2L:55:5MU:C6	2.52	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:4E:80:ILE:HG23	5:4E:91:LEU:HD23	1.99	0.45
33:69:41:GLU:HG3	33:69:41:GLU:H	1.56	0.45
2:12:111:ARG:HA	2:12:111:ARG:HD3	1.80	0.45
26:1H:25:U:H2'	26:1H:26:G:C8	2.52	0.45
1:1G:257:G:H1	1:1G:269:C:H42	1.66	0.45
26:14:2022:U:O2'	26:14:2617:C:H5'	2.17	0.45
1:1G:129(A):G:C2	1:1G:191(A):G:C8	3.05	0.45
33:61:5:LEU:HD13	33:61:13:GLY:O	2.17	0.45
1:13:861:G:H4'	8:7E:18:ARG:HH21	1.81	0.45
29:21:134:ILE:O	29:21:137:HIS:HB2	2.17	0.45
1:13:317:G:C6	1:13:318:G:C6	3.05	0.45
38:55:13:HIS:HD2	38:55:15:SER:H	1.65	0.45
1:13:895:G:H2'	1:13:896:C:C6	2.52	0.45
7:62:101:LEU:O	7:62:105:VAL:HG23	2.17	0.45
22:1K:44:G:H1'	22:1K:45:U:C5	2.52	0.45
35:25:71:ARG:NH2	35:25:122:LEU:O	2.48	0.45
31:49:111:LEU:HA	31:49:111:LEU:HD23	1.82	0.45
1:1G:881:G:C6	1:1G:882:C:C4	3.05	0.45
47:E5:72:ARG:HB3	47:E5:75:LEU:HB2	1.98	0.45
1:1G:1523:G:OP1	11:2A:123:LYS:HD2	2.16	0.45
36:35:135:LEU:HD13	36:35:139:LYS:HE2	1.99	0.45
44:F8:41:ASN:O	44:F8:45:THR:HG23	2.17	0.45
26:14:1045:A:H1'	26:14:1047:G:C4	2.52	0.44
26:1H:1371:G:H2'	26:1H:1372:U:H5	1.82	0.44
41:85:92:ARG:CZ	42:95:11:GLN:H	2.29	0.44
1:13:1002:G:C4	1:13:1003:G:C8	3.05	0.44
1:1G:1289:A:OP1	21:1B:10:ARG:NE	2.50	0.44
21:1B:3:LYS:O	21:1B:14:TRP:CE3	2.70	0.44
1:1G:983:A:H61	1:1G:1222:G:H22	1.64	0.44
30:31:63:LYS:HG2	30:31:65:TRP:O	2.17	0.44
36:35:144:GLU:HA	36:35:145:PRO:HD3	1.74	0.44
4:3E:22:LYS:HB2	4:3E:26:CYS:SG	2.57	0.44
1:1G:554:C:H2'	1:1G:555:C:C6	2.52	0.44
26:14:638:G:H2'	26:14:639:U:O4'	2.16	0.44
29:29:11:MET:HG3	29:29:24:THR:N	2.30	0.44
26:14:947:G:N2	26:14:971:C:C2	2.85	0.44
33:69:76:THR:HG21	33:69:139:GLN:O	2.17	0.44
1:1G:222:U:C2	1:1G:223:U:C5	3.06	0.44
1:1G:45:U:H2'	1:1G:46:G:H8	1.81	0.44
8:72:29:SER:H	8:72:32:LYS:HB2	1.81	0.44
46:H8:62:PRO:O	46:H8:63:ASP:HB2	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
30:39:107:LYS:HD3	30:39:107:LYS:HA	1.76	0.44
26:1H:1348:G:H2'	26:1H:1349:A:H5''	1.98	0.44
26:14:2773:C:OP1	29:29:166:THR:OG1	2.36	0.44
26:1H:1519:G:O2'	26:1H:1520:U:H5'	2.16	0.44
43:A5:79:GLY:CA	43:A5:100:THR:HG22	2.47	0.44
39:A8:67:ARG:HB2	39:A8:67:ARG:CZ	2.48	0.44
26:14:1858:G:H2'	26:14:1883:G:H22	1.81	0.44
45:G8:43:ASN:OD1	45:G8:65:ALA:HB3	2.17	0.44
26:1H:2793:G:H8	26:1H:2793:G:OP2	1.99	0.44
31:49:33:ARG:NH2	31:49:162:THR:HG21	2.31	0.44
26:1H:532:A:N7	26:1H:2021:C:O2'	2.35	0.44
26:14:1652:A:OP1	38:55:8:ARG:NH1	2.50	0.44
44:F8:21:PHE:O	44:F8:23:GLU:O	2.33	0.44
8:7E:28:ALA:HB3	8:7E:57:PRO:HB2	1.99	0.44
1:13:1151:A:H5'	10:1I:41:PRO:HA	1.99	0.44
26:1H:937:U:H2'	26:1H:938:G:O4'	2.17	0.44
26:1H:664:C:H4'	26:1H:941:A:OP1	2.17	0.44
1:1G:123:C:OP1	1:1G:312:C:H5'	2.17	0.44
1:13:338:A:C6	1:13:339:C:C4	3.05	0.44
26:14:1510:A:H2'	26:14:1511:A:O4'	2.17	0.44
30:39:31:HIS:O	30:39:31:HIS:CG	2.70	0.44
9:8E:23:ASN:HD22	9:8E:23:ASN:H	1.64	0.44
31:49:18:GLU:O	31:49:21:ARG:HB3	2.16	0.44
26:1H:2744:G:N2	32:51:143:GLN:OE1	2.49	0.44
7:62:78:ARG:HB2	7:62:156:TRP:CZ3	2.52	0.44
51:M8:38:LYS:HA	51:M8:40:HIS:H	1.82	0.44
26:1H:2291:U:O2'	26:1H:2374:C:O2	2.27	0.44
9:82:77:ILE:O	9:82:81:ILE:HG12	2.17	0.44
26:1H:1386:C:OP2	26:1H:1396:U:H5	2.01	0.44
22:1K:46:7MG:H5''	22:1K:46:7MG:H82	1.98	0.44
1:13:739:C:C4	1:13:740:U:C5	3.05	0.44
36:78:125:VAL:O	36:78:144:GLU:HB2	2.17	0.44
27:1J:55:U:O2'	27:1J:56:G:H5'	2.17	0.44
39:A8:87:PHE:CE2	39:A8:89:ARG:HB2	2.52	0.44
26:14:1702:G:H2'	26:14:1703:G:O4'	2.17	0.44
26:14:2472:G:N2	26:14:2477:C:OP2	2.47	0.44
26:14:2688:U:C5	26:14:2720:U:OP2	2.69	0.44
24:3K:39:U:H1'	24:3K:40:C:C6	2.52	0.44
49:G5:47:ASN:C	49:G5:49:LYS:H	2.17	0.44
1:1G:848:C:H2'	1:1G:849:C:H6	1.82	0.44
24:3L:4:C:H2'	24:3L:5:G:O4'	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:1H:30:G:H2'	26:1H:31:C:C6	2.53	0.44
39:A8:29:PHE:CD1	39:A8:30:ARG:N	2.85	0.44
13:4I:108:ARG:N	13:4I:108:ARG:HD2	2.32	0.44
42:D8:43:GLU:HA	42:D8:44:LYS:HA	1.70	0.44
32:51:10:PRO:HG2	32:51:50:VAL:O	2.17	0.44
1:1G:1329:A:H5'	13:4A:25:ILE:O	2.17	0.44
28:19:12:SER:O	28:19:16:MET:HB2	2.18	0.44
3:22:58:GLU:O	3:22:59:ARG:HG3	2.17	0.44
39:65:3:ARG:NH2	39:65:4:LEU:HB2	2.33	0.44
26:1H:236:C:H2'	26:1H:237:C:C6	2.50	0.44
42:D8:8:GLY:O	42:D8:10:LYS:HE3	2.17	0.44
21:1F:12:LYS:HB3	21:1F:22:ARG:HD2	1.98	0.44
1:1G:160:A:H2'	1:1G:161:A:O4'	2.17	0.44
37:88:104:PHE:O	37:88:105:GLU:HB3	2.17	0.44
48:F5:41:ARG:HD3	48:F5:43:TYR:CE1	2.51	0.44
45:G8:30:VAL:O	45:G8:32:PRO:HD3	2.17	0.44
6:52:11:ASN:OD1	6:52:12:PRO:HD2	2.17	0.44
1:13:1417:G:C6	1:13:1482:G:C6	3.06	0.44
26:1H:2692:C:O2'	26:1H:2693:A:H5'	2.17	0.44
11:2I:80:VAL:HG22	11:2I:103:LEU:HD12	1.99	0.44
26:14:299:A:OP2	58:14:3888:HOH:O	2.21	0.44
26:14:483:A:H4'	45:C5:49:VAL:HA	2.00	0.44
35:68:34:THR:OG1	35:68:35:VAL:N	2.50	0.44
26:14:1344:G:H4'	26:14:1384:A:C5	2.52	0.44
1:13:1499:A:O2'	1:13:1520:G:H5'	2.16	0.44
1:1G:940:C:H2'	1:1G:941:G:C8	2.53	0.44
11:2A:20:TYR:HB2	11:2A:31:THR:HG23	1.98	0.44
43:A5:19:LEU:HB3	52:J5:25:LEU:HD12	2.00	0.44
29:21:182:LEU:HD12	29:21:183:LEU:H	1.82	0.44
4:3E:117:ALA:O	4:3E:120:LEU:HB2	2.17	0.44
36:35:65:ARG:HB3	36:35:65:ARG:NH1	2.31	0.44
41:C8:79:PHE:C	41:C8:79:PHE:CD2	2.91	0.44
4:3E:24:GLU:HG2	4:3E:24:GLU:H	1.43	0.44
41:C8:16:LYS:HE2	41:C8:16:LYS:HB3	1.72	0.44
1:13:1272:G:C6	1:13:1273:G:C5	3.05	0.44
10:1A:89:ASP:HB3	10:1A:91:PRO:HD3	1.98	0.44
31:41:13:GLU:O	31:41:14:GLU:HB3	2.16	0.44
26:14:2657:A:O2'	32:59:160:LYS:HE3	2.16	0.44
26:1H:2153:G:O6	26:1H:2154:G:N2	2.50	0.44
1:1G:533:A:H2'	58:1G:1728:HOH:O	2.17	0.44
26:1H:2445:G:OP1	30:31:74:ARG:NH2	2.51	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1G:1362:C:H2'	1:1G:1362(A):C:H5''	1.99	0.44
36:35:19:VAL:HG13	36:35:21:ARG:N	2.21	0.44
36:35:39:LYS:HB2	36:35:45:LEU:HD11	1.99	0.44
26:14:993:G:C6	26:14:994:C:N4	2.85	0.44
26:14:849:A:H3'	26:14:850:C:C6	2.52	0.44
26:14:1786:A:H1'	26:14:1938:A:N6	2.32	0.44
26:1H:1814:G:N1	58:1H:4231:HOH:O	2.50	0.44
26:14:2286:A:C8	26:14:2287:A:C6	3.05	0.44
26:14:2287:A:N6	26:14:2344:U:H3	2.08	0.44
31:41:37:VAL:HG23	31:41:99:MET:CE	2.47	0.44
1:1G:545:C:OP2	4:32:65:ARG:NH2	2.51	0.44
35:68:68:GLU:HA	35:68:78:ARG:HB3	1.99	0.44
1:1G:1118:C:H1'	1:1G:1179:A:C4	2.52	0.44
13:4A:84:ILE:C	13:4A:86:CYS:H	2.20	0.44
26:14:2298:A:N6	26:14:2318:G:C8	2.85	0.44
2:1E:70:PHE:HE2	2:1E:90:MET:CB	2.31	0.44
4:32:108:LEU:HD23	4:32:110:PHE:HE2	1.81	0.44
1:13:667:G:H4'	15:6I:51:HIS:ND1	2.32	0.44
26:1H:582:G:H2'	26:1H:583:G:H8	1.82	0.44
13:4A:96:LEU:HD22	13:4A:103:THR:HG21	1.99	0.44
47:I8:19:LYS:HD3	47:I8:19:LYS:HA	1.65	0.44
33:69:3:VAL:HA	33:69:39:ALA:HB2	1.99	0.44
26:1H:1449:A:H5'	26:1H:1449(A):G:OP2	2.16	0.44
26:1H:1060:U:O2	26:1H:1088:A:H8	2.01	0.44
28:11:159:ALA:HB1	28:11:198:ASN:O	2.18	0.44
7:62:69:VAL:HG13	7:62:134:ALA:O	2.17	0.44
47:I8:28:GLY:N	47:I8:67:VAL:O	2.42	0.44
3:22:18:TRP:HE1	14:5A:55:GLY:H	1.65	0.44
26:14:1260:G:C6	26:14:1261:C:C4	3.05	0.44
26:1H:44:A:C2'	26:1H:45:G:H5'	2.47	0.44
26:1H:2246:G:H2'	26:1H:2247:A:H8	1.81	0.44
26:1H:2863:C:H2'	26:1H:2864:G:H8	1.83	0.44
1:13:464:G:C6	1:13:466:C:H5'	2.52	0.44
26:14:297:C:H2'	26:14:298:G:O4'	2.18	0.44
26:1H:1092:C:H2'	26:1H:1093:G:H5'	1.99	0.44
37:45:118:LEU:HD12	37:45:131:ILE:HG23	1.99	0.44
1:13:490:G:H2'	1:13:491:G:C8	2.52	0.44
41:C8:28:ARG:HD3	41:C8:38:THR:OG1	2.17	0.44
1:1G:64:G:H4'	1:1G:65:U:O5'	2.17	0.44
10:1I:47:PHE:CZ	14:5I:37:PHE:HE1	2.35	0.44
4:3E:61:LYS:HG3	4:3E:203:VAL:HG13	1.98	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:2453:A:H2'	26:14:2454:G:O4'	2.18	0.44
2:1E:72:GLY:HA2	2:1E:165:VAL:HG22	1.99	0.44
38:98:59:ASP:OD1	38:98:59:ASP:N	2.36	0.44
12:3I:127:GLU:N	12:3I:127:GLU:OE1	2.50	0.44
1:1G:1186:G:OP2	1:1G:1186:G:H8	2.00	0.44
26:1H:2547:U:H2'	26:1H:2548:G:C8	2.52	0.44
26:1H:2273:A:H2'	26:1H:2274:A:C8	2.52	0.44
26:14:531:C:OP1	26:14:561:G:N2	2.51	0.44
15:6I:17:ARG:HG2	15:6I:21:ASP:OD2	2.17	0.44
2:12:75:LYS:HA	2:12:78:GLN:CB	2.40	0.44
27:16:11:C:O5'	27:16:12:C:H5	2.01	0.44
31:41:66:GLN:NE2	31:41:93:THR:O	2.42	0.44
1:13:1029:G:O2'	1:13:1032(A):G:N2	2.51	0.44
41:C8:92:ARG:HH11	41:C8:95:LEU:CD2	2.31	0.44
1:1G:191:G:C6	1:1G:192:U:C4	3.05	0.44
26:14:2250:G:C2	37:45:82:ARG:HB3	2.52	0.44
29:21:120:TRP:CD2	29:21:155:LYS:HG2	2.53	0.44
1:13:607:A:H2	16:7I:31:LYS:HG3	1.82	0.44
34:58:96:GLU:HG2	34:58:97:ARG:H	1.82	0.44
30:31:127:GLU:HG2	30:31:196:LEU:HD12	2.00	0.44
26:1H:2699:C:H2'	26:1H:2700:C:O4'	2.18	0.44
5:42:90:VAL:HG23	5:42:121:LYS:O	2.18	0.44
5:4E:110:LEU:HD13	5:4E:118:ILE:HG21	1.97	0.44
53:O8:21:TYR:HB2	53:O8:22:ALA:H	1.43	0.44
5:4E:37:ARG:HA	5:4E:113:ALA:HA	1.99	0.44
1:1G:116:A:O5'	1:1G:116:A:H8	2.00	0.44
26:1H:270(L):U:H3	33:6I:50:ARG:CG	2.30	0.44
7:62:94:ARG:O	7:62:97:GLN:HB3	2.18	0.44
26:14:303:U:H2'	26:14:304:G:H8	1.81	0.44
34:15:56:ASN:HD22	34:15:126:PRO:HA	1.83	0.44
20:BI:97:ALA:O	20:BI:99:LEU:N	2.51	0.44
35:25:87:ILE:HG23	35:25:88:ASN:O	2.18	0.44
1:13:1218:C:OP2	14:5I:9:LYS:NZ	2.49	0.44
50:L8:32:GLN:NE2	50:L8:32:GLN:HA	2.32	0.44
26:1H:839:U:O2'	26:1H:1191:G:N3	2.50	0.44
23:2L:33:OMC:HM22	23:2L:34:U:H5'	1.98	0.44
27:16:32:C:C2	27:16:51:G:N2	2.86	0.44
1:13:690:G:H2'	1:13:691:G:O4'	2.16	0.44
37:88:66:ILE:O	37:88:104:PHE:N	2.51	0.44
10:1A:76:ASN:HB3	10:1A:78:ASN:HD22	1.83	0.44
26:1H:2430:A:H8	26:1H:2431:U:H5	1.65	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1G:1254:C:OP1	10:1A:45:ARG:HA	2.17	0.44
49:K8:33:MET:HG2	49:K8:37:PHE:CE1	2.52	0.44
26:1H:8:A:H2'	26:1H:9:U:C6	2.52	0.44
26:1H:2705:A:O2'	26:1H:2852:G:OP1	2.23	0.44
26:14:673:C:H4'	30:39:82:ILE:HG12	1.99	0.44
11:2I:73:MET:HG2	11:2I:103:LEU:HD23	1.99	0.44
26:14:270(G):C:H2'	26:14:270(H):C:H6	1.81	0.44
50:H5:46:ASN:O	50:H5:50:VAL:HG22	2.17	0.44
1:13:279:A:C5	17:8I:98:LEU:HD12	2.53	0.44
44:B5:5:TYR:CE1	49:G5:30:ARG:HG3	2.52	0.44
26:1H:2296:U:H4'	26:1H:2297:C:OP1	2.17	0.44
26:1H:1370:C:HO2'	26:1H:1811:G:HO2'	1.66	0.44
7:6E:72:ARG:HG3	7:6E:142:GLU:OE1	2.18	0.44
1:13:595:G:N2	1:13:643:C:N4	2.65	0.44
26:14:675:A:C4	26:14:804:A:C2	3.05	0.44
10:1I:26:ALA:C	10:1I:30:SER:HB3	2.37	0.44
1:1G:935:A:H2'	1:1G:936:C:C6	2.52	0.44
17:8I:28:PRO:HA	17:8I:34:LYS:O	2.16	0.44
26:14:777:A:C2	26:14:778:G:C4	3.05	0.44
26:1H:1750:G:C2	26:1H:1751:C:C5	3.05	0.44
26:14:873:G:N2	26:14:905:U:C2	2.85	0.44
26:14:1973:G:H2'	26:14:1974:C:C6	2.53	0.44
32:59:144:VAL:O	32:59:148:ILE:HG12	2.18	0.44
1:13:1336:C:H5''	1:13:1336:C:C6	2.52	0.44
38:98:28:LEU:HA	38:98:28:LEU:HD23	1.88	0.44
24:3K:34:G:OP1	24:3K:34:G:H8	2.00	0.44
34:58:16:ILE:HB	34:58:54:VAL:HG22	1.99	0.44
26:1H:1065:U:H2'	26:1H:1066:U:O4'	2.18	0.44
1:1G:1127:G:H1'	1:1G:1148:U:N3	2.32	0.44
26:1H:568:U:OP1	36:78:36:LYS:HE3	2.18	0.44
26:1H:2582:G:C2	26:1H:2583:G:C8	3.06	0.44
26:14:2134:A:H2	26:14:2159:G:O2'	2.01	0.44
19:AA:40:ILE:HA	19:AA:44:MET:SD	2.57	0.44
53:O8:41:PRO:C	53:O8:43:CYS:H	2.21	0.44
29:21:119:ARG:CG	29:21:119:ARG:HH11	2.28	0.44
8:7E:87:SER:HB2	8:7E:93:VAL:N	2.32	0.44
1:1G:836:G:C6	1:1G:851:G:C6	3.05	0.44
26:1H:1050:A:C8	26:1H:2751:G:C8	3.06	0.44
1:1G:1302:U:C5	13:4A:17:VAL:HG21	2.53	0.44
29:29:102:VAL:HB	29:29:199:ARG:O	2.17	0.44
26:1H:1024:G:C3'	26:1H:1025:G:H5''	2.46	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1G:115:G:H1'	1:1G:116:A:N7	2.33	0.44
26:14:1013:C:N3	26:14:1149:G:N2	2.44	0.44
4:3E:98:GLU:O	4:3E:103:ASN:ND2	2.50	0.44
26:1H:1055:G:H1'	26:1H:1085:A:H2	1.81	0.44
26:1H:1102:C:H2'	26:1H:1103:A:H8	1.81	0.44
1:1G:1227:A:O2'	13:4A:115:LYS:HE2	2.18	0.44
28:11:36:PRO:HA	28:11:61:LEU:HD12	2.00	0.44
29:21:82:ARG:O	29:21:84:PHE:N	2.50	0.44
1:1G:183:G:H1	1:1G:194:C:H42	1.65	0.44
14:5I:6:LEU:HB3	14:5I:23:ARG:NH2	2.32	0.44
1:1G:994:A:C2	14:5A:5:ALA:HB2	2.53	0.44
1:1G:762:C:H2'	1:1G:763:G:C8	2.52	0.44
34:15:120:LEU:HG	34:15:122:VAL:HG23	1.98	0.44
26:14:610:C:H2'	26:14:611:C:C6	2.52	0.44
51:M8:54:GLY:HA2	51:M8:57:GLU:HB3	1.99	0.44
1:13:115:G:C2	1:13:289:G:N7	2.86	0.44
26:1H:270(G):C:H2'	26:1H:270(H):C:O4'	2.18	0.44
26:14:1432:C:H2'	26:14:1433:U:O4'	2.17	0.44
43:E8:85:VAL:HA	43:E8:95:ILE:HG22	2.00	0.44
29:21:11:MET:HG2	29:21:24:THR:HA	1.99	0.44
36:78:23:PRO:O	36:78:25:SER:N	2.50	0.44
5:4E:48:ALA:HB3	5:4E:54:ALA:HB2	2.00	0.44
29:29:170:LEU:HA	29:29:170:LEU:HD13	1.58	0.44
16:7I:82:GLN:H	16:7I:82:GLN:HG2	1.67	0.44
12:3I:85:ILE:HA	12:3I:85:ILE:HD13	1.73	0.44
47:I8:36:ILE:O	47:I8:36:ILE:HD13	2.17	0.44
30:39:144:LYS:HA	30:39:144:LYS:HD2	1.52	0.44
54:P8:8:ASN:C	54:P8:8:ASN:OD1	2.56	0.44
20:BI:71:THR:HG22	20:BI:72:LEU:N	2.24	0.44
26:14:572:A:H2'	26:14:573:G:O4'	2.18	0.44
31:41:98:ARG:HH21	51:M8:1:MET:HG3	1.82	0.44
1:1G:157:G:H1	1:1G:164:U:H3	1.64	0.44
26:14:362:U:H3'	26:14:363:G:H5''	1.99	0.44
26:14:2494:G:C5	26:14:2495:G:N7	2.86	0.44
47:I8:48:GLY:N	47:I8:79:VAL:O	2.43	0.44
16:7I:77:ALA:HB3	16:7I:79:VAL:HG23	2.00	0.44
30:31:32:LEU:CD2	30:31:105:VAL:HG13	2.44	0.44
1:1G:1376:U:OP1	7:62:98:SER:OG	2.26	0.44
9:82:10:ARG:HH21	9:82:11:LYS:HG3	1.83	0.44
4:3E:173:TRP:HA	4:3E:187:ARG:HG2	2.00	0.44
23:2L:19:G:H4'	23:2L:20:G:OP1	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:3E:110:PHE:CE2	4:3E:148:VAL:HG23	2.52	0.44
45:C5:57:GLN:HB3	45:C5:58:GLY:H	1.53	0.44
26:14:234:C:H2'	26:14:235:U:H6	1.82	0.44
26:14:2:G:N2	26:14:2900:A:H61	2.15	0.44
7:62:91:VAL:HG12	7:62:95:ARG:HB3	1.99	0.44
26:1H:2430:A:H8	26:1H:2431:U:C5	2.35	0.44
17:8I:78:GLU:HG2	17:8I:81:ARG:HD2	1.99	0.44
9:8E:10:ARG:HD2	9:8E:11:LYS:HG3	2.00	0.44
29:29:37:ARG:HA	29:29:42:ASP:OD1	2.17	0.44
20:BA:26:ASN:HB3	20:BA:71:THR:OG1	2.18	0.44
1:1G:321:A:O2'	1:1G:322:C:H5'	2.18	0.44
33:61:118:LYS:HA	33:61:119:PRO:HD3	1.73	0.44
1:13:825:G:C6	1:13:826:C:C4	3.05	0.44
26:14:1449(A):G:H2'	26:14:1450:C:C6	2.53	0.44
1:1G:149:A:H2'	1:1G:150:C:H6	1.83	0.44
7:62:78:ARG:HB2	7:62:156:TRP:HZ3	1.83	0.44
44:B5:5:TYR:HB3	49:G5:33:MET:HB2	1.99	0.44
13:4I:105:THR:OG1	13:4I:106:ASN:N	2.51	0.44
1:1G:818:G:O2'	1:1G:819:A:H5'	2.18	0.44
28:11:119:ALA:CB	28:11:130:ALA:HB3	2.47	0.44
1:1G:1455:G:H5'	20:BA:32:ALA:HB2	2.00	0.44
10:1A:42:THR:HG23	10:1A:67:THR:O	2.17	0.44
8:72:42:GLU:HG3	8:72:109:ILE:HD12	1.99	0.44
26:1H:216:A:H2'	26:1H:217:G:H8	1.82	0.44
28:11:109:ASP:HB2	28:11:197:GLY:HA3	2.00	0.44
55:Q8:26:LYS:O	55:Q8:28:GLY:N	2.50	0.44
49:G5:29:LYS:HG2	49:G5:57:ILE:HD13	1.99	0.44
26:14:1225:C:H4'	42:95:85:LYS:CG	2.47	0.44
26:1H:2428:G:N2	36:78:61:ARG:NH2	2.66	0.44
27:1J:66:A:C6	27:1J:107:U:H2'	2.53	0.44
30:39:187:VAL:HG13	36:35:1:MET:O	2.18	0.44
12:3A:27:LEU:HD11	12:3A:62:SER:OG	2.17	0.44
10:1I:54:PHE:HB3	10:1I:55:LYS:H	1.70	0.44
27:16:42:C:C6	31:41:69:ALA:HB2	2.53	0.44
26:14:1826:G:H2'	26:14:1827:C:C6	2.53	0.44
53:O8:45:LYS:HA	53:O8:45:LYS:HD3	1.66	0.44
7:6E:71:PRO:HA	7:6E:138:LYS:HG2	1.99	0.44
1:13:1284:C:H2'	1:13:1285:A:N7	2.32	0.44
1:13:1226:C:O3'	13:4I:111:LYS:NZ	2.49	0.44
13:4A:30:ALA:O	13:4A:34:LEU:HG	2.18	0.44
36:78:52:GLU:CG	36:78:57:THR:HA	2.47	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1G:1325:C:H5''	21:1B:17:THR:HG21	1.98	0.44
27:16:15:A:H1'	27:16:109:G:C4	2.51	0.44
44:B5:18:TYR:C	44:B5:20:GLY:N	2.69	0.44
26:1H:2099:U:H2'	26:1H:2100:G:C8	2.53	0.44
46:D5:99:TYR:HA	46:D5:124:ILE:O	2.17	0.44
1:13:1327:C:P	21:1F:12:LYS:HZ1	2.41	0.44
26:14:231:C:C2'	26:14:232:G:H5'	2.48	0.44
26:14:2257:U:O4	58:14:3651:HOH:O	2.21	0.44
17:8A:59:ILE:HD13	17:8A:73:VAL:HA	1.99	0.44
4:3E:116:GLN:NE2	4:3E:157:LEU:HD11	2.32	0.44
6:52:15:ASP:O	6:52:19:LEU:HB2	2.18	0.44
54:P8:10:ARG:HD3	54:P8:14:LYS:HD3	2.00	0.44
26:14:1805:U:C2'	26:14:1806:C:H5'	2.48	0.44
27:16:3:C:H2'	27:16:4:C:H6	1.82	0.44
27:16:3:C:H2'	27:16:4:C:C6	2.52	0.44
8:7E:84:ARG:O	8:7E:135:CYS:HB2	2.17	0.44
1:13:1402:C:H2'	1:13:1403:C:O4'	2.17	0.44
1:1G:453:A:H4'	16:7A:72:ARG:HB2	1.98	0.44
1:1G:646:U:H2'	1:1G:647:C:C6	2.52	0.44
2:12:32:ILE:HD12	2:12:41:ILE:O	2.18	0.44
2:12:128:GLU:O	2:12:130:ARG:HG2	2.18	0.44
26:1H:671:C:OP1	36:78:42:SER:O	2.36	0.44
1:1G:236:G:H2'	1:1G:237:C:O4'	2.17	0.44
1:13:256:U:H2'	1:13:257:G:C8	2.53	0.44
26:14:1157:G:C2	26:14:1158:C:C2	3.06	0.44
43:A5:36:LEU:HD11	43:A5:47:VAL:HG12	1.99	0.44
1:13:900:A:H8	1:13:900:A:O5'	2.01	0.44
32:51:149:ARG:HG2	32:51:149:ARG:O	2.17	0.44
26:1H:1225:C:O2'	42:D8:85:LYS:HA	2.18	0.44
26:1H:1965:C:H3'	26:1H:1966:A:H2'	1.99	0.44
4:32:175:SER:HB3	4:32:186:LEU:HD11	1.98	0.44
26:1H:1771:C:C1'	26:1H:1786:A:C8	3.00	0.44
26:1H:1786:A:H1'	26:1H:1938:A:N6	2.33	0.44
36:78:17:LYS:HE2	36:78:27:HIS:NE2	2.32	0.44
51:M8:39:CYS:H	51:M8:41:PRO:HD3	1.83	0.44
2:12:6:THR:HG21	2:12:217:ARG:HB3	2.00	0.44
26:1H:49:A:C8	26:1H:120:U:H5	2.34	0.44
14:5A:23:ARG:HA	14:5A:23:ARG:HD3	1.79	0.44
34:15:67:LEU:O	34:15:88:GLU:HG3	2.17	0.44
11:2A:57:THR:HG22	11:2A:58:PRO:HD2	1.98	0.44
26:14:2312:U:O2	31:49:42:GLY:HA3	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
27:1J:2:C:H2'	27:1J:3:C:C5	2.53	0.44
28:19:53:PHE:O	28:19:218:ARG:HB2	2.18	0.44
26:14:827:U:H2'	26:14:2430:A:H2	1.82	0.44
48:J8:73:LEU:HD11	48:J8:95:LEU:HD21	1.99	0.44
26:1H:1466:G:N3	26:1H:1547:C:N4	2.66	0.44
26:14:94:G:O2'	49:G5:46:GLN:HB3	2.18	0.44
36:35:62:LEU:HD13	36:35:63:PRO:HD2	2.00	0.44
26:14:2295:C:OP1	39:65:10:ARG:NH1	2.46	0.44
16:7I:28:ARG:NH1	16:7I:29:ASP:OD1	2.39	0.44
40:75:36:GLU:HG3	40:75:36:GLU:O	2.18	0.44
26:14:1278:A:O3'	38:55:34:ILE:HG13	2.17	0.44
26:14:2611:U:O2'	52:J5:3:LYS:HG2	2.18	0.44
3:22:33:LEU:O	3:22:36:ASP:N	2.51	0.44
1:1G:624:C:H2'	1:1G:625:G:C8	2.53	0.44
18:9A:22:VAL:HG12	18:9A:55:ARG:O	2.17	0.44
26:1H:2400:G:H4'	53:O8:18:ARG:O	2.18	0.44
28:11:44:ASN:O	28:11:46:GLN:O	2.35	0.44
26:14:184:C:H2'	26:14:185:U:C6	2.53	0.44
34:58:43:THR:HA	34:58:44:PRO:HD2	1.79	0.44
1:13:458:C:H42	1:13:475:G:H1	1.65	0.44
3:22:199:LYS:HB3	3:22:201:TYR:HE1	1.82	0.44
34:58:128:HIS:HB2	34:58:129:PRO:HD2	1.99	0.44
28:19:166:GLN:HE22	28:19:176:ARG:NH2	2.15	0.44
31:49:120:LEU:HD22	31:49:133:LEU:HD11	1.99	0.44
37:88:17:LEU:HB3	37:88:39:PRO:HB2	1.99	0.44
1:1G:1232:U:H2'	1:1G:1233:G:O4'	2.17	0.44
26:1H:276:A:C5	26:1H:278:A:H2	2.36	0.44
1:1G:323:U:O4'	20:BA:19:SER:HB2	2.18	0.44
26:1H:557:U:H2'	26:1H:558:G:H8	1.83	0.44
26:14:699:A:H2'	26:14:700:G:O4'	2.17	0.44
48:F5:90:ILE:HG22	48:F5:91:LYS:N	2.33	0.44
14:5A:38:GLY:O	14:5A:39:LEU:HD23	2.18	0.44
26:14:2461:C:H2'	26:14:2462:U:C6	2.52	0.44
26:14:55:G:C2	26:14:116:C:C2	3.05	0.44
12:3A:110:VAL:CG2	12:3A:120:TYR:HB3	2.48	0.44
26:14:1101:U:H2'	26:14:1102:C:H6	1.83	0.44
26:14:843:G:N2	26:14:936:C:C2	2.86	0.44
1:13:255:G:P	17:8I:69:LYS:HZ3	2.40	0.44
1:1G:1115:C:H1'	14:5A:61:TRP:HB2	1.99	0.44
10:1I:38:ILE:HG23	10:1I:71:LEU:HB3	1.99	0.44
32:51:46:GLU:OE1	32:51:51:ARG:NH2	2.51	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:1695:G:OP1	28:19:7:LYS:NZ	2.51	0.44
1:13:27:G:H4'	4:3E:209:ARG:HG3	1.99	0.44
23:2K:2:G:N3	23:2K:2:G:H2'	2.32	0.44
29:21:63:LEU:O	29:21:63:LEU:HD23	2.17	0.44
1:1G:575:G:H4'	1:1G:575:G:OP1	2.18	0.44
15:6A:3:ILE:H	15:6A:3:ILE:HG12	1.47	0.44
28:11:273:ARG:HG2	28:11:273:ARG:HH21	1.82	0.44
26:14:2666:C:H3'	26:14:2667:C:H6	1.83	0.44
26:1H:2820:A:O2'	26:1H:2821:A:OP1	2.35	0.44
1:13:1348:U:H2'	1:13:1349:A:H8	1.82	0.44
12:3I:113:ARG:O	12:3I:114:LYS:HD2	2.17	0.44
26:14:996:A:H4'	41:85:92:ARG:HD3	1.98	0.44
26:1H:70:G:H21	26:1H:71:A:N6	2.16	0.44
26:14:2402:C:H5	26:14:2415:G:H22	1.65	0.44
24:3K:3:C:H2'	24:3K:4:C:O4'	2.18	0.44
48:J8:91:LYS:CG	48:J8:92:LYS:H	2.30	0.44
31:41:67:LYS:HE3	31:41:67:LYS:O	2.18	0.44
16:7I:74:LEU:HD22	16:7I:79:VAL:HG21	2.00	0.44
1:1G:1321:C:H4'	13:4A:87:TYR:CE1	2.53	0.44
28:19:71:ASP:CG	28:19:103:ARG:HH12	2.21	0.44
1:13:278:G:N2	17:8I:95:TYR:HB3	2.33	0.44
26:1H:1675:C:H2'	26:1H:1676:A:O4'	2.17	0.44
1:1G:1347:G:N2	1:1G:1373:G:H2'	2.33	0.44
1:1G:1028(A):C:C4	1:1G:1028(B):C:H5	2.36	0.44
1:1G:1028:C:N4	1:1G:1033:G:H1	2.13	0.44
26:1H:287:C:H2'	26:1H:288:C:C6	2.48	0.44
35:25:104:ARG:HB3	35:25:104:ARG:CZ	2.48	0.44
26:14:1785:A:H4'	26:14:1982:C:O2'	2.18	0.44
7:62:26:PHE:O	7:62:30:ILE:HG13	2.18	0.44
9:8E:89:ASN:O	9:8E:91:ASP:N	2.50	0.44
52:N8:42:PRO:O	52:N8:44:THR:HG22	2.18	0.44
55:Q8:34:TRP:HD1	55:Q8:34:TRP:C	2.19	0.44
29:29:62:PRO:C	29:29:64:LYS:N	2.71	0.44
26:1H:932:G:H4'	26:1H:933:A:O5'	2.17	0.44
26:14:26:G:C6	26:14:27:G:N1	2.85	0.44
1:1G:338:A:H2'	1:1G:339:C:C6	2.53	0.44
26:14:2694:G:C5	26:14:2695:C:C5	3.06	0.44
1:1G:944:G:C2	1:1G:1340:A:C6	3.06	0.44
44:B5:12:VAL:HG13	44:B5:27:THR:OG1	2.18	0.44
1:1G:243:A:C2	1:1G:245:C:C2	3.05	0.44
1:13:1279:A:O2'	1:13:1281:U:OP2	2.26	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
18:9A:66:LEU:O	18:9A:70:ILE:HG13	2.18	0.44
26:14:1636:C:H2'	26:14:1637:A:C8	2.53	0.44
27:16:89:G:H2'	27:16:89(A):A:C8	2.53	0.44
1:1G:66:G:C2	1:1G:67:C:C6	3.05	0.44
37:88:17:LEU:HA	37:88:17:LEU:HD23	1.58	0.44
26:14:2648:C:H2'	26:14:2649:U:O4'	2.18	0.44
15:6I:82:ILE:O	15:6I:86:GLY:N	2.50	0.44
44:B5:5:TYR:HD1	49:G5:33:MET:SD	2.41	0.44
1:1G:272:C:H2'	1:1G:273:A:C8	2.53	0.44
29:21:67:PHE:C	29:21:69:LYS:H	2.21	0.44
26:1H:2670:A:O2'	26:1H:2671:A:H5'	2.18	0.44
26:1H:39:C:H2'	26:1H:40:C:C6	2.52	0.44
34:15:116:LEU:O	34:15:119:ARG:N	2.38	0.44
30:31:80:ALA:HA	30:31:81:PRO:HD3	1.88	0.44
26:14:1405:U:H2'	26:14:1406:U:C6	2.53	0.44
1:13:1338:G:C6	1:13:1339:A:C6	3.05	0.44
1:1G:302:G:H4'	12:3A:17:LYS:HE2	2.00	0.44
26:14:1894:C:O2'	26:14:1895:C:H5'	2.17	0.44
6:5E:25:ILE:HD13	6:5E:25:ILE:HA	1.83	0.44
26:1H:2256:G:N2	26:1H:2275:C:N4	2.66	0.43
24:3K:8:U:H2'	24:3K:13:C:H5	1.81	0.43
46:H8:97:GLU:HB3	46:H8:125:LEU:HD11	2.00	0.43
26:1H:444:C:O2'	26:1H:445:C:H5'	2.17	0.43
27:1J:2:C:H2'	27:1J:3:C:C6	2.53	0.43
1:1G:363:A:C5	12:3A:31:PRO:HD2	2.53	0.43
48:J8:73:LEU:HD21	48:J8:95:LEU:HD23	1.99	0.43
32:51:4:ILE:O	32:51:4:ILE:HG12	2.17	0.43
26:1H:1063:G:H22	26:1H:1076:C:C2'	2.31	0.43
1:13:658:G:H2'	1:13:659:U:H6	1.82	0.43
26:1H:1668:A:OP1	35:68:5:GLN:HG3	2.18	0.43
26:14:2016:U:H2'	26:14:2017:U:C6	2.52	0.43
38:98:34:ILE:HG22	38:98:114:VAL:HB	1.99	0.43
33:61:120:ILE:HG12	33:61:126:TYR:CE2	2.53	0.43
1:1G:539:A:H2'	1:1G:540:G:C8	2.53	0.43
26:1H:1025:G:H8	26:1H:1025:G:OP1	1.99	0.43
31:49:53:LEU:CD2	31:49:87:PRO:HB2	2.48	0.43
14:5A:17:LYS:NZ	14:5A:18:VAL:HG13	2.33	0.43
26:14:2611:U:O2'	26:14:2612:C:O5'	2.30	0.43
26:1H:910:A:N1	26:1H:2277:G:H1'	2.32	0.43
15:6I:39:LEU:O	15:6I:42:HIS:N	2.50	0.43
8:7E:8:ASP:OD2	8:7E:12:ARG:HD2	2.17	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:13:1289:A:H3'	1:13:1290:G:C8	2.52	0.43
19:AA:66:MET:H	19:AA:67:VAL:HB	1.82	0.43
45:G8:63:LYS:HG3	45:G8:64:GLU:H	1.83	0.43
26:14:2773:C:H2'	26:14:2774:C:H6	1.83	0.43
27:1J:99:A:C4	27:1J:100:G:C8	3.06	0.43
26:1H:950:G:C5	26:1H:951:C:C4	3.06	0.43
45:G8:6:HIS:N	45:G8:6:HIS:CD2	2.86	0.43
26:14:2784:C:H1'	29:29:37:ARG:NH2	2.33	0.43
31:49:60:LEU:HD21	31:49:92:VAL:HG11	1.99	0.43
5:4E:43:LEU:HD23	5:4E:133:TYR:HD1	1.83	0.43
26:14:511:U:C5	26:14:512:G:C5	3.05	0.43
26:1H:431:U:O2'	26:1H:432:A:H5'	2.18	0.43
26:14:1374:G:H2'	26:14:1375:C:C6	2.53	0.43
53:K5:12:GLU:HA	53:K5:23:THR:HB	2.00	0.43
17:8A:10:VAL:HG21	17:8A:52:LYS:O	2.18	0.43
26:1H:2028:U:H2'	26:1H:2029:G:O4'	2.17	0.43
44:B5:26:TYR:CE2	44:B5:89:ILE:HG13	2.53	0.43
28:11:148:GLU:HB2	28:11:151:LYS:HD2	2.00	0.43
26:1H:1474:C:H2'	26:1H:1475:G:C8	2.53	0.43
1:13:752:G:H4'	15:6I:69:TYR:OH	2.18	0.43
26:1H:1754:C:P	40:B8:96:ARG:HH12	2.40	0.43
20:BA:87:LYS:HE3	20:BA:91:LEU:HD11	2.00	0.43
5:4E:6:PHE:HD1	5:4E:6:PHE:HA	1.74	0.43
26:1H:1153:C:C4	26:1H:1154:G:C5	3.06	0.43
46:H8:67:LEU:HD22	46:H8:90:VAL:HG11	2.00	0.43
7:6E:115:ARG:HB3	7:6E:118:VAL:HG12	1.98	0.43
1:1G:976:G:N2	1:1G:1362:C:H2'	2.33	0.43
26:14:2552:U:C2	26:14:2554:U:H5''	2.53	0.43
26:1H:577:G:OP1	26:1H:2502:G:O2'	2.30	0.43
1:13:61:G:H2'	1:13:62:U:O4'	2.18	0.43
31:41:61:ALA:HA	31:41:66:GLN:O	2.18	0.43
46:H8:14:LYS:HA	46:H8:15:PRO:HD2	1.72	0.43
26:1H:1532:C:H42	26:1H:1539:G:H1	1.67	0.43
26:14:878:A:H2'	26:14:879:G:O4'	2.17	0.43
26:1H:1465:G:C5	26:1H:1466:G:N7	2.86	0.43
28:19:41:GLY:HA3	28:19:43:ARG:HD3	2.00	0.43
1:1G:1014:A:H4'	19:AA:14:HIS:HE1	1.79	0.43
1:1G:1347:G:N2	1:1G:1374:A:OP2	2.41	0.43
39:A8:84:GLN:OE1	39:A8:110:LEU:HD13	2.19	0.43
26:14:1639:U:H5'	58:14:3552:HOH:O	2.17	0.43
48:F5:92:LYS:O	48:F5:93:GLU:C	2.54	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1G:828:A:C4	1:1G:859:A:C8	3.06	0.43
1:13:429:U:H1'	1:13:430:A:H5''	1.98	0.43
26:14:1149:G:H2'	26:14:1150:C:C6	2.53	0.43
9:8E:27:THR:O	9:8E:63:ILE:N	2.49	0.43
37:88:43:THR:HG22	37:88:94:VAL:HG12	1.99	0.43
29:29:9:VAL:HA	40:75:3:ARG:HG3	2.00	0.43
26:1H:322:A:H5'	26:1H:340:A:C1'	2.48	0.43
45:G8:39:VAL:O	45:G8:42:VAL:HG22	2.16	0.43
31:41:47:LYS:NZ	31:41:81:LYS:HB2	2.32	0.43
10:1A:45:ARG:HB3	10:1A:65:LEU:HB3	1.99	0.43
34:58:28:THR:HG22	34:58:29:LYS:N	2.33	0.43
1:1G:523:A:H61	12:3A:92:ASP:HB2	1.83	0.43
45:C5:48:ALA:HB1	45:C5:50:ARG:HD3	2.00	0.43
1:13:45:U:H2'	1:13:46:G:C8	2.53	0.43
26:14:2619:C:H2'	26:14:2620:C:C6	2.52	0.43
26:14:2512:C:H2'	26:14:2513:G:O4'	2.17	0.43
26:1H:39:C:O2	30:31:46:ARG:NH2	2.51	0.43
30:31:46:ARG:HA	30:31:46:ARG:HD2	1.70	0.43
26:14:822:U:O2'	26:14:823:G:H5'	2.18	0.43
1:13:1486:G:H2'	1:13:1487:G:O4'	2.18	0.43
26:1H:2019:A:C6	26:1H:2020:A:N7	2.86	0.43
26:1H:2320:A:H2'	26:1H:2320:A:N3	2.32	0.43
1:1G:1496:C:H2'	1:1G:1497:G:C8	2.53	0.43
36:78:79:ARG:HB2	36:78:110:TYR:HD1	1.82	0.43
20:BI:92:LEU:O	20:BI:96:GLY:HA3	2.19	0.43
1:13:869:G:O5'	1:13:869:G:H8	2.01	0.43
26:1H:2259:G:C2	26:1H:2282:G:C6	3.06	0.43
1:1G:1072:G:C6	1:1G:1073:U:C4	3.07	0.43
39:65:38:GLN:HG2	39:65:47:THR:HG21	1.98	0.43
9:8E:29:ASN:OD1	9:8E:65:VAL:N	2.51	0.43
26:1H:1387:C:O2	26:1H:1388:G:C8	2.71	0.43
1:1G:476:G:H2'	1:1G:477:G:H8	1.83	0.43
26:1H:1389:G:C2	26:1H:1399:C:O2	2.72	0.43
1:13:963:G:N2	10:1I:55:LYS:NZ	2.66	0.43
26:14:1771:C:H1'	26:14:1786:A:C8	2.53	0.43
1:13:376:G:O3'	16:7I:5:ARG:HD2	2.19	0.43
1:1G:1048:G:N2	1:1G:1209:C:N3	2.62	0.43
26:14:2336:A:H61	47:E5:43:THR:HB	1.83	0.43
2:12:204:ASN:N	2:12:204:ASN:OD1	2.52	0.43
26:14:1946:U:H2'	26:14:1947:C:C6	2.54	0.43
26:1H:1049:C:H1'	26:1H:1113:U:H4'	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:78:24:GLY:O	36:78:26:GLY:N	2.51	0.43
13:4A:54:VAL:HA	13:4A:57:ARG:HB3	1.99	0.43
29:21:35:GLN:HB3	29:21:48:GLN:HB2	2.00	0.43
1:13:1320:C:H2'	1:13:1321:C:O4'	2.18	0.43
1:13:1320:C:H42	19:AI:36:ARG:HG3	1.83	0.43
29:21:13:ARG:HG2	29:21:13:ARG:NH1	2.33	0.43
26:14:1151:G:H4'	41:85:81:HIS:CG	2.53	0.43
26:14:2153:G:HO2'	26:14:2154:G:H8	1.63	0.43
26:14:2893:G:H4'	26:14:2894:G:O5'	2.18	0.43
26:1H:1085:A:H1'	26:1H:1086:A:C2	2.53	0.43
26:14:1795:C:H2'	26:14:1796:U:H6	1.83	0.43
53:K5:18:ARG:HH12	53:K5:43:CYS:HB2	1.83	0.43
26:14:140:A:H8	26:14:1408:C:O2'	1.98	0.43
3:22:21:ARG:O	3:22:58:GLU:HA	2.18	0.43
3:22:40:ARG:HA	3:22:43:LEU:HB2	2.00	0.43
20:BI:56:MET:HB3	20:BI:56:MET:HE2	1.88	0.43
29:21:65:GLY:HA2	29:21:70:ALA:CB	2.49	0.43
26:1H:2144:U:H1'	26:1H:2148:G:N2	2.32	0.43
32:51:86:GLU:HG3	32:51:165:ALA:H	1.82	0.43
26:1H:2330:G:H2'	26:1H:2331:G:O4'	2.17	0.43
1:1G:1105:A:H2'	1:1G:1106:G:C8	2.52	0.43
11:2I:57:THR:HA	11:2I:58:PRO:HD2	1.71	0.43
26:14:1298:C:C2'	26:14:1302:A:HO2'	2.24	0.43
26:14:2567:G:H2'	26:14:2568:C:H6	1.83	0.43
49:G5:38:GLN:O	49:G5:43:GLN:HA	2.18	0.43
26:14:1511:A:H2'	26:14:1512:G:O4'	2.19	0.43
1:13:595:G:H22	1:13:643:C:N4	2.16	0.43
30:39:64:ILE:HG13	30:39:65:TRP:CE2	2.53	0.43
26:1H:270(X):G:C6	26:1H:270(Y):G:N1	2.87	0.43
35:68:31:LYS:HB3	35:68:32:TYR:CE2	2.54	0.43
26:14:390:A:C6	36:35:71:VAL:HG11	2.53	0.43
40:75:57:PHE:HA	40:75:79:HIS:CD2	2.53	0.43
33:69:4:ILE:HG21	33:69:47:LEU:HD13	2.00	0.43
1:13:150:C:H2'	1:13:151:A:O4'	2.18	0.43
26:14:142:G:H5''	26:14:1598:C:O2'	2.17	0.43
26:14:1152:C:H4'	41:85:77:SER:HA	2.00	0.43
1:1G:800:G:H8	1:1G:800:G:O5'	2.01	0.43
8:7E:39:LEU:HB3	8:7E:45:ILE:HG12	2.01	0.43
30:31:155:LEU:HB2	30:31:189:THR:HG21	2.00	0.43
38:55:103:ARG:CZ	38:55:110:PRO:HD3	2.48	0.43
7:6E:111:ARG:HB3	7:6E:111:ARG:HH11	1.83	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:13:536:C:H2'	1:13:537:G:H8	1.81	0.43
5:42:107:ARG:CZ	5:42:107:ARG:HB3	2.48	0.43
26:14:996:A:N3	26:14:997:G:C8	2.86	0.43
55:Q8:48:PHE:CE2	55:Q8:52:LYS:HB2	2.54	0.43
26:1H:1782:C:H2'	26:1H:2608:G:O2'	2.19	0.43
26:1H:1395:A:P	58:1H:3645:HOH:O	2.70	0.43
2:1E:21:ARG:C	2:1E:23:ARG:H	2.22	0.43
26:14:1060:U:H5'	26:14:1061:U:C6	2.53	0.43
29:21:188:VAL:HA	29:21:189:PRO:HD3	1.77	0.43
11:2A:85:ARG:HE	11:2A:111:ASP:HB3	1.82	0.43
1:13:660:G:H2'	1:13:661:G:C8	2.53	0.43
26:14:890:A:H2'	26:14:892:G:C8	2.50	0.43
4:3E:18:LYS:HE3	4:3E:18:LYS:HB2	1.63	0.43
4:32:199:ASN:CB	4:32:202:LEU:HG	2.48	0.43
1:1G:843:U:H3'	1:1G:848:C:O4'	2.18	0.43
39:A8:84:GLN:O	39:A8:85:VAL:HG13	2.18	0.43
1:1G:1273:G:H3'	1:1G:1274:G:H8	1.83	0.43
27:1J:113:C:O2'	39:65:46:VAL:HG13	2.18	0.43
1:1G:862:C:H2'	1:1G:863:U:H5'	2.01	0.43
1:13:1226:C:H4'	1:13:1227:A:OP1	2.19	0.43
26:1H:2314:C:H2'	26:1H:2315:G:C8	2.50	0.43
26:14:2893:G:H5'	26:14:2894:G:OP1	2.18	0.43
26:1H:1086:A:H1'	26:1H:1103:A:N6	2.33	0.43
26:14:2542:A:H1'	26:14:2543:G:C8	2.53	0.43
14:5I:4:LYS:O	14:5I:7:ILE:HG13	2.19	0.43
17:8A:63:ARG:HG2	17:8A:64:PRO:CD	2.48	0.43
32:59:76:VAL:O	32:59:80:SER:OG	2.20	0.43
20:BA:81:LYS:O	20:BA:85:MET:HG2	2.18	0.43
26:1H:303:U:H2'	26:1H:304:G:C8	2.52	0.43
37:88:87:LYS:HA	37:88:87:LYS:HD2	1.47	0.43
26:14:492:A:H2'	26:14:493:G:O4'	2.17	0.43
26:14:296:C:H2'	26:14:297:C:H6	1.83	0.43
1:1G:1084:G:H2'	1:1G:1085:U:H6	1.84	0.43
31:49:10:LYS:O	31:49:15:VAL:HG23	2.17	0.43
1:13:651:C:H2'	1:13:652:U:C6	2.54	0.43
8:72:118:VAL:O	8:72:119:LEU:HD23	2.19	0.43
26:1H:1195:G:N3	26:1H:1227:A:H2	2.16	0.43
26:1H:1471:A:C2	26:1H:1472:A:C8	3.06	0.43
26:14:564:C:H5''	42:95:75:PHE:CZ	2.54	0.43
14:5I:13:THR:HG23	14:5I:20:ALA:HB2	2.00	0.43
2:1E:25:ASN:HA	2:1E:26:PRO:HD3	1.84	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:500:G:N2	26:14:502:A:H3'	2.33	0.43
48:J8:81:LYS:HD2	48:J8:81:LYS:N	2.34	0.43
26:1H:572:A:H5''	26:1H:573:G:OP2	2.17	0.43
26:14:1416:G:O2'	26:14:1417:C:O4'	2.35	0.43
36:78:19:VAL:HG21	36:78:27:HIS:CG	2.54	0.43
22:1K:14:A:C6	22:1K:22:G:N3	2.86	0.43
40:B8:91:ARG:O	40:B8:116:ALA:HA	2.19	0.43
4:32:25:ARG:HG2	4:32:30:LYS:O	2.18	0.43
50:L8:7:LYS:HD2	50:L8:34:GLU:HG2	1.99	0.43
16:7I:3:LYS:O	16:7I:21:VAL:HA	2.18	0.43
8:7E:34:GLU:HB3	8:7E:118:VAL:HG21	2.01	0.43
49:K8:51:ARG:NH1	49:K8:55:ARG:HH12	2.16	0.43
29:29:33:VAL:HG12	29:29:89:ASP:CB	2.47	0.43
31:41:37:VAL:H	31:41:99:MET:HE3	1.82	0.43
46:D5:53:ILE:HA	46:D5:70:LEU:HD22	2.01	0.43
26:14:249:C:H5''	58:14:3522:HOH:O	2.18	0.43
26:1H:287:C:O2'	26:1H:288:C:H5'	2.19	0.43
26:1H:2262:U:H4'	26:1H:2328:A:C2	2.54	0.43
7:62:92:SER:O	7:62:96:GLN:HG3	2.19	0.43
1:1G:980:C:H3'	1:1G:981:U:H6	1.79	0.43
28:19:44:ASN:HD21	28:19:46:GLN:HG3	1.82	0.43
26:14:990:A:H5'	26:14:990:A:C8	2.49	0.43
26:1H:65:C:H2'	26:1H:66:C:C6	2.54	0.43
26:1H:720:C:H2'	26:1H:721:C:C6	2.54	0.43
35:68:71:ARG:NH2	35:68:77:ILE:HG21	2.33	0.43
6:52:60:PHE:C	6:52:61:LEU:HD12	2.39	0.43
20:BA:11:SER:HA	20:BA:13:LEU:HB2	1.99	0.43
45:C5:52:SER:OG	45:C5:56:PRO:HA	2.18	0.43
1:13:1448:C:H42	1:13:1455:G:H1	1.67	0.43
26:14:1317:A:H2'	26:14:1318:C:H6	1.83	0.43
26:1H:2181:G:H2'	26:1H:2182:G:C8	2.53	0.43
1:1G:791:G:C6	1:1G:792:A:N7	2.87	0.43
15:6A:75:PRO:HA	15:6A:78:TYR:HB3	1.99	0.43
26:1H:2846:G:H2'	26:1H:2847:U:O4'	2.18	0.43
1:13:1507:A:OP2	25:4K:13:A:N6	2.52	0.43
26:1H:2862:G:H2'	26:1H:2863:C:H6	1.83	0.43
26:14:1858:G:H8	26:14:1858:G:OP2	2.02	0.43
26:14:2086:U:H2'	26:14:2087:G:C8	2.54	0.43
45:G8:89:PHE:CD1	45:G8:90:LEU:N	2.87	0.43
14:5A:59:ALA:HB1	14:5A:61:TRP:HZ3	1.83	0.43
30:31:81:PRO:HB3	30:31:89:VAL:HG23	1.99	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:1H:2208:U:H4'	28:11:151:LYS:HG2	2.00	0.43
26:14:286:C:H2'	26:14:287:C:C6	2.53	0.43
26:14:2815:C:H5'	52:J5:29:THR:HG21	2.00	0.43
38:55:81:ASP:O	38:55:82:GLU:HB3	2.18	0.43
26:1H:1820:U:H4'	26:1H:1821:A:OP2	2.19	0.43
26:14:117:G:C6	26:14:119:A:C6	3.07	0.43
26:14:1213:A:H2'	26:14:1214:A:C8	2.53	0.43
26:1H:478:A:C6	26:1H:480:A:C6	3.06	0.43
30:31:34:TRP:CZ2	36:78:8:PRO:HB3	2.53	0.43
35:25:89:ASN:N	35:25:89:ASN:OD1	2.52	0.43
26:14:1088:A:H2'	26:14:1088:A:N3	2.32	0.43
16:7I:25:ARG:HG3	16:7I:25:ARG:HH11	1.84	0.43
53:K5:45:LYS:N	53:K5:45:LYS:HD2	2.33	0.43
18:9I:21:LYS:HD2	18:9I:21:LYS:HA	1.86	0.43
17:8I:101:ARG:NH2	17:8I:101:ARG:HB2	2.32	0.43
22:1K:60:U:H5'	22:1K:61:C:OP2	2.18	0.43
26:1H:2572:A:N7	29:21:144:ARG:HD2	2.34	0.43
26:1H:1387:C:C2	26:1H:1388:G:C8	3.05	0.43
26:14:1225:C:O3'	42:95:85:LYS:HA	2.18	0.43
26:1H:1264:G:H3'	26:1H:1265:A:H5''	1.99	0.43
26:1H:2127:G:H2'	26:1H:2128:C:O4'	2.18	0.43
31:49:96:ARG:C	31:49:98:ARG:H	2.22	0.43
33:69:130:TYR:C	33:69:131:LYS:HD2	2.38	0.43
4:32:4:TYR:HE2	4:32:11:LEU:HD11	1.83	0.43
26:1H:2303:G:O2'	31:41:132:ASN:HB2	2.18	0.43
10:1A:61:GLU:OE1	14:5A:58:LYS:HE2	2.18	0.43
26:14:2395:C:H2'	26:14:2396:G:O4'	2.18	0.43
39:65:10:ARG:O	39:65:14:VAL:HG22	2.18	0.43
13:4A:7:VAL:HG21	31:49:115:ARG:NE	2.34	0.43
26:14:270(S):G:H2'	26:14:270(T):G:H8	1.83	0.43
1:1G:711:G:O2'	1:1G:712:A:H5'	2.18	0.43
1:1G:373:A:C2	1:1G:374:A:C8	3.07	0.43
26:14:162:U:H4'	26:14:171:G:O4'	2.18	0.43
13:4A:25:ILE:O	13:4A:29:ARG:HB2	2.19	0.43
23:2K:64:G:C2	23:2K:65:G:N7	2.87	0.43
41:C8:36:ARG:HD3	41:C8:40:PHE:CZ	2.54	0.43
32:51:130:ARG:HH11	32:51:130:ARG:HB3	1.84	0.43
1:13:280:C:O2	17:8I:38:ARG:HG3	2.19	0.43
26:1H:1568:G:P	28:11:63:ARG:HH12	2.42	0.43
26:14:2543:G:H1'	26:14:2766:G:H5'	2.01	0.43
26:1H:1108:U:C4	26:1H:1109:C:N4	2.87	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
40:75:86:ILE:HG21	40:75:86:ILE:HD13	1.82	0.43
5:42:127:ASN:HA	5:42:128:PRO:HD3	1.91	0.43
1:1G:829:G:O2'	1:1G:830:G:H5'	2.18	0.43
32:51:121:ILE:HG12	32:51:140:LYS:HD3	2.00	0.43
26:14:511:U:O4	26:14:512:G:N1	2.50	0.43
1:13:416:G:H2'	1:13:417:C:C6	2.54	0.43
49:G5:33:MET:O	49:G5:37:PHE:HD1	2.02	0.43
26:14:557:U:H2'	26:14:558:G:C8	2.54	0.43
26:1H:803:U:C4	26:1H:804:A:N7	2.86	0.43
8:7E:23:SER:HA	8:7E:61:VAL:O	2.19	0.43
26:14:2408:U:H2'	26:14:2409:G:C8	2.54	0.43
1:1G:1472:U:H2'	1:1G:1473:A:O4'	2.19	0.43
26:1H:1922:G:H2'	26:1H:1923:U:C6	2.54	0.43
28:19:108:PRO:HB3	28:19:143:HIS:HE2	1.83	0.43
26:1H:2712:U:H1'	26:1H:2712(A):A:C8	2.53	0.43
27:16:73:A:C4	27:16:104:A:C2	3.07	0.43
33:61:29:TYR:O	33:61:32:PRO:HD2	2.18	0.43
1:1G:583:A:H2'	1:1G:584:G:O4'	2.19	0.43
26:14:1233:C:H2'	26:14:1234:U:H6	1.84	0.43
30:31:184:TYR:CE1	36:78:3:LEU:HD11	2.53	0.43
47:E5:50:ASN:O	47:E5:62:LEU:HB2	2.18	0.43
6:5E:46:ARG:HG3	6:5E:47:ARG:N	2.34	0.43
1:13:609:A:H2'	1:13:610:G:H5'	2.01	0.43
3:22:140:ARG:NE	3:22:140:ARG:HA	2.34	0.43
42:95:66:ARG:H	42:95:66:ARG:HG2	1.68	0.43
26:1H:315:G:C6	26:1H:316:C:C4	3.06	0.43
26:1H:851:U:O2'	50:L8:42:ALA:O	2.37	0.43
48:J8:85:LEU:HA	48:J8:85:LEU:HD13	1.72	0.43
45:G8:85:VAL:O	45:G8:86:ARG:HD3	2.19	0.43
16:7A:40:ASP:HB3	16:7A:48:TRP:CB	2.39	0.43
1:1G:241:C:C2	1:1G:286:G:C2	3.07	0.43
20:BI:26:ASN:HB3	20:BI:71:THR:OG1	2.19	0.43
3:22:148:GLY:O	3:22:203:PHE:HB3	2.19	0.43
26:1H:1389:G:C2	26:1H:1390:U:O2	2.72	0.43
26:1H:2680:C:O3'	29:21:8:LYS:NZ	2.52	0.43
1:1G:1368:G:OP2	9:82:112:LYS:HD2	2.18	0.43
1:13:265:G:H5''	17:8I:65:ILE:O	2.19	0.43
1:13:101:A:C4	1:13:102:G:C8	3.06	0.43
26:1H:1332:G:N2	26:1H:1610:A:N7	2.67	0.43
26:14:1425:G:H2'	26:14:1426:G:C8	2.53	0.43
1:13:1104:G:H4'	2:1E:111:ARG:NH1	2.34	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:1H:286:C:H2'	26:1H:287:C:C6	2.53	0.43
26:1H:550:G:O2'	26:1H:1220:A:N3	2.38	0.43
32:51:12:PRO:HB3	32:51:48:GLY:HA2	2.00	0.43
37:88:32:TYR:CE1	37:88:133:ARG:HG3	2.54	0.43
26:14:2346:A:C2	26:14:2383:G:C2	3.06	0.43
34:58:133:GLN:CD	34:58:133:GLN:H	2.22	0.43
14:5A:12:ARG:HG3	14:5A:12:ARG:H	1.38	0.43
2:12:208:ILE:HA	2:12:211:ILE:HD12	1.99	0.43
26:1H:234:C:H2'	26:1H:235:U:C6	2.50	0.43
1:1G:316:G:H2'	1:1G:317:G:C8	2.54	0.43
26:14:1176:G:H5'	26:14:1177:A:OP1	2.19	0.43
50:L8:8:LEU:HD13	50:L8:31:LEU:HA	2.00	0.43
5:42:92:LYS:HE2	8:72:105:ARG:NH2	2.34	0.43
1:13:1422:G:C5'	35:68:48:PRO:HB3	2.49	0.43
10:1A:32:ALA:HA	10:1A:76:ASN:HB2	2.01	0.43
26:1H:2335:A:N7	26:1H:2337:G:C5	2.87	0.43
1:13:240:C:H2'	1:13:241:C:H6	1.82	0.43
11:2I:59:TYR:CE2	11:2I:63:LEU:HD11	2.53	0.43
30:31:178:PRO:HB2	30:31:201:VAL:CG2	2.48	0.43
26:1H:2280:G:H2'	26:1H:2281:C:H5'	2.00	0.43
1:13:617:G:H5'	16:7I:45:THR:HG23	2.00	0.43
30:39:126:VAL:HG13	30:39:193:VAL:HG23	2.01	0.43
46:H8:4:ARG:HD3	46:H8:60:GLU:OE2	2.18	0.43
46:H8:4:ARG:HB3	46:H8:58:VAL:HG23	1.99	0.43
1:1G:141:A:H1'	1:1G:182:U:O2	2.19	0.43
26:1H:4:C:H42	26:1H:2899:G:H1	1.65	0.43
32:59:117:PRO:HA	32:59:118:PRO:HD2	1.88	0.43
33:69:135:GLU:OE2	33:69:135:GLU:N	2.50	0.43
50:L8:6:VAL:HG11	50:L8:47:VAL:HG22	2.01	0.43
52:N8:16:ARG:HG3	52:N8:17:ASP:N	2.34	0.43
26:1H:2467:C:H4'	37:88:123:HIS:CG	2.53	0.43
24:3L:71:G:H2'	24:3L:72:C:H5''	2.01	0.43
26:1H:1779:U:H2'	58:1H:3625:HOH:O	2.18	0.43
1:1G:1148:U:OP1	9:82:7:THR:HG21	2.19	0.43
42:95:85:LYS:HG3	42:95:87:HIS:CA	2.48	0.43
40:B8:88:ILE:O	40:B8:88:ILE:HG13	2.18	0.43
26:14:563:G:H5'	26:14:572:A:H4'	2.01	0.43
27:16:11:C:H3'	27:16:12:C:H6	1.82	0.43
1:1G:1352:C:OP1	21:1B:3:LYS:NZ	2.38	0.43
3:22:91:LEU:O	3:22:95:THR:OG1	2.32	0.43
4:32:8:VAL:HA	4:32:11:LEU:HD12	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:78:121:LYS:HB3	36:78:121:LYS:HE2	1.82	0.43
9:82:111:ARG:HG2	9:82:112:LYS:H	1.84	0.43
1:1G:1512:U:H2'	1:1G:1513:A:C8	2.53	0.43
34:15:15:LEU:HD23	34:15:134:ARG:HD2	2.00	0.43
28:11:260:ARG:HH22	28:11:266:SER:HB3	1.83	0.43
1:13:410:G:H5''	1:13:411:A:OP1	2.18	0.43
13:4I:2:ALA:O	13:4I:10:PRO:HD2	2.18	0.43
13:4I:108:ARG:NH1	13:4I:111:LYS:HB3	2.34	0.43
40:75:107:ASP:N	40:75:107:ASP:OD1	2.51	0.43
31:49:56:ALA:HB2	31:49:153:ARG:NH2	2.34	0.43
36:35:110:TYR:HB3	36:35:111:ARG:H	1.72	0.43
12:3I:32:PHE:HB3	12:3I:84:LEU:HD11	2.01	0.43
17:8I:10:VAL:HG13	17:8I:19:VAL:HB	2.01	0.43
29:29:173:VAL:N	29:29:183:LEU:O	2.37	0.43
26:14:1171:G:O2'	26:14:1173:G:O4'	2.20	0.43
19:AA:66:MET:HA	19:AA:67:VAL:O	2.18	0.43
12:3A:75:HIS:HA	12:3A:111:LYS:NZ	2.34	0.43
5:42:92:LYS:HG2	5:42:93:PRO:HD2	2.00	0.43
3:2E:148:GLY:HA3	3:2E:172:ARG:O	2.19	0.43
55:M5:32:LEU:HA	55:M5:32:LEU:HD12	1.63	0.43
26:14:2239:G:H5'	28:19:251:GLY:HA3	2.01	0.43
1:1G:109:A:H5'	1:1G:110:C:H5	1.84	0.43
26:14:2340:G:H2'	26:14:2341:G:C8	2.54	0.43
26:1H:2840:C:H42	26:1H:2877:G:H1	1.67	0.43
26:14:826:U:H2'	26:14:828:U:O4'	2.19	0.43
26:1H:270(R):G:H2'	26:1H:270(S):G:C8	2.53	0.43
31:49:55:LYS:HG2	31:49:150:ASP:OD1	2.19	0.43
53:O8:33:LYS:O	53:O8:35:GLU:HG3	2.19	0.43
21:1B:6:ARG:HH11	21:1B:15:ARG:NH2	2.16	0.43
32:59:56:SER:OG	32:59:57:ASP:N	2.52	0.43
2:1E:135:GLN:O	2:1E:139:LYS:HB2	2.17	0.43
12:3A:10:LEU:HD23	12:3A:10:LEU:HA	1.86	0.43
26:1H:1954:G:O2'	26:1H:1956:U:O4	2.30	0.43
26:14:957:A:N6	26:14:2459:A:C8	2.87	0.43
37:88:2:LEU:HD12	37:88:2:LEU:HA	1.75	0.43
1:13:13:U:O2	1:13:914:A:H3'	2.18	0.43
1:1G:1350:A:C2	1:1G:1351:U:C2	3.06	0.43
46:H8:18:LEU:O	46:H8:21:ALA:HB3	2.18	0.43
1:1G:651:C:H2'	1:1G:652:U:O4'	2.18	0.43
1:13:1375:A:P	7:6E:28:ASN:HD22	2.42	0.43
1:1G:537:G:H5''	12:3A:113:ARG:HH12	1.80	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:2751:G:N1	32:59:2:SER:HB3	2.34	0.43
47:E5:20:ARG:HD3	47:E5:20:ARG:HA	1.76	0.43
26:14:831:G:H5''	26:14:832:G:OP2	2.19	0.43
26:1H:2130:U:H1'	26:1H:2159:G:N2	2.33	0.43
26:1H:72:U:OP2	44:F8:1:MET:N	2.51	0.43
1:1G:674:G:H2'	1:1G:675:A:C8	2.54	0.43
26:14:2685:G:P	40:75:51:ARG:HH12	2.41	0.43
26:1H:1063:G:N1	26:1H:1076:C:O2	2.51	0.43
8:7E:87:SER:HB2	8:7E:93:VAL:H	1.83	0.43
28:19:242:ARG:N	28:19:242:ARG:HD3	2.32	0.43
1:13:266:G:H5''	1:13:267:C:C5	2.53	0.43
28:19:41:GLY:C	28:19:43:ARG:H	2.22	0.43
28:19:49:ILE:HD11	28:19:52:ARG:HA	2.00	0.43
46:D5:155:LEU:HD12	46:D5:163:LEU:HD13	2.00	0.43
26:1H:1313:U:H4'	26:1H:1332:G:H4'	2.00	0.43
46:D5:70:LEU:O	46:D5:89:PHE:N	2.46	0.43
1:1G:1028(A):C:H42	1:1G:1032(B):G:N2	2.16	0.43
13:4I:13:LYS:O	13:4I:44:ARG:NH2	2.51	0.43
1:13:992:U:O2	1:13:993:G:N2	2.52	0.43
26:1H:1799:G:H5'	26:1H:1819:A:N6	2.33	0.43
26:1H:507:A:C5'	26:1H:508:G:H5'	2.47	0.43
26:14:629:G:H1	26:14:634:C:N4	2.14	0.43
26:14:2298:A:N6	26:14:2318:G:H2'	2.34	0.43
26:1H:849:A:H5''	26:1H:850:C:OP2	2.18	0.43
26:1H:548:A:C5	26:1H:549:G:H1'	2.53	0.43
26:1H:2858:C:H2'	26:1H:2859:G:O4'	2.19	0.43
1:1G:755:G:OP2	15:6A:65:ARG:HD3	2.18	0.43
4:3E:92:VAL:O	4:3E:96:LEU:HD22	2.19	0.43
28:11:17:THR:HG22	28:11:205:VAL:H	1.82	0.43
39:65:54:LEU:O	39:65:56:LEU:N	2.46	0.43
50:L8:8:LEU:HB2	50:L8:28:LEU:HD22	2.00	0.43
29:21:60:ASN:HB2	29:21:61:ARG:H	1.62	0.43
22:1K:69:G:H3'	22:1K:70:G:H5''	2.01	0.43
3:2E:134:ILE:HD11	3:2E:153:VAL:HG21	2.00	0.43
26:1H:1110:G:O2'	26:1H:1111:A:C8	2.72	0.43
26:14:1599:C:C2	26:14:1600:C:C5	3.06	0.43
54:L5:16:HIS:HB2	54:L5:44:PRO:HG2	2.00	0.43
26:14:1693:U:H4'	26:14:1694:C:OP2	2.19	0.43
26:1H:208:C:H2'	26:1H:209:C:C6	2.53	0.43
1:1G:949:A:C2	1:1G:1233:G:N3	2.87	0.43
1:13:945:G:C2	1:13:946:A:C8	3.06	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
42:95:7:THR:HG23	42:95:22:VAL:HG21	2.00	0.43
26:1H:2259:G:C2	26:1H:2282:G:N1	2.86	0.43
26:1H:2288:A:C2	26:1H:2325:G:C8	3.07	0.43
31:41:28:VAL:O	31:41:31:VAL:HG13	2.19	0.43
9:82:2:GLU:HG2	9:82:3:GLN:HG3	2.00	0.43
26:1H:1279:G:N2	26:1H:1292:U:C2	2.87	0.43
13:4A:112:GLY:HA3	13:4A:113:PRO:HD3	1.78	0.43
1:13:864:A:H5''	1:13:865:A:OP2	2.19	0.43
26:14:1348:G:H5''	26:14:1349:A:OP2	2.19	0.43
26:14:254:G:O6	55:M5:5:LYS:HG2	2.19	0.43
26:14:725:G:H8	26:14:725:G:O5'	2.02	0.43
28:19:112:GLN:O	28:19:115:GLN:HG3	2.18	0.43
2:1E:180:LEU:O	2:1E:181:PHE:HB2	2.19	0.43
26:14:270(V):G:H2'	26:14:270(W):G:O4'	2.19	0.43
27:1J:13:A:N1	27:1J:69:G:O2'	2.45	0.43
26:14:997:G:P	41:85:58:ARG:HH21	2.41	0.43
46:H8:72:ARG:NH1	46:H8:89:PHE:HD2	2.17	0.43
55:Q8:14:VAL:HG11	55:Q8:21:LYS:NZ	2.33	0.43
5:4E:126:ARG:CG	5:4E:126:ARG:HH11	2.28	0.43
19:AA:52:TYR:HB2	19:AA:57:HIS:CE1	2.54	0.43
29:29:16:ARG:O	29:29:17:ASP:HB2	2.19	0.43
38:55:38:VAL:CG1	38:55:42:LYS:HD2	2.45	0.43
8:72:68:ARG:NH1	8:72:70:GLN:HG2	2.34	0.43
1:13:1126:U:N3	1:13:1127:G:C2	2.87	0.43
29:29:119:ARG:HA	29:29:160:TYR:CE2	2.54	0.43
30:39:7:TYR:O	30:39:15:SER:HA	2.19	0.43
26:14:1570:A:H2'	26:14:1571:A:C8	2.54	0.43
1:13:636:U:H2'	1:13:637:G:C8	2.53	0.43
26:14:1778:U:P	58:14:3501:HOH:O	2.77	0.43
26:14:2378:A:O3'	39:65:23:ARG:HD2	2.19	0.43
47:E5:26:TYR:HB2	47:E5:29:GLN:OE1	2.19	0.43
24:3L:9:A:H61	24:3L:22:G:H5'	1.84	0.43
26:14:2637:U:C2	26:14:2782:G:N2	2.87	0.43
20:BI:53:LEU:O	20:BI:57:ARG:NH1	2.52	0.43
39:65:56:LEU:HB3	39:65:58:LEU:HD22	2.01	0.43
4:3E:108:LEU:HD23	4:3E:110:PHE:CE1	2.54	0.43
1:13:1096:C:H2'	1:13:1097:C:C6	2.50	0.43
26:1H:1448:G:N2	26:1H:1449:A:N6	2.66	0.43
26:14:2228:G:C5	26:14:2229:C:C4	3.07	0.43
1:1G:277:C:H5''	17:8A:68:ARG:NH2	2.34	0.43
32:51:92:ILE:HG13	32:51:92:ILE:H	1.48	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1G:1396:A:H4'	1:1G:1397:C:OP2	2.19	0.43
46:H8:53:ILE:HA	46:H8:71:VAL:HG13	2.00	0.43
33:69:44:LEU:HD23	33:69:44:LEU:HA	1.85	0.43
26:14:977:G:H2'	26:14:978:G:H8	1.83	0.43
2:1E:51:LEU:HD23	2:1E:201:ILE:HD12	2.01	0.43
26:14:128:C:H2'	26:14:129:C:H6	1.84	0.43
53:O8:27:LYS:H	53:O8:27:LYS:CE	2.31	0.43
8:7E:82:HIS:HE1	8:7E:136:GLU:OE2	2.02	0.43
1:13:721:G:C6	1:13:733:A:C2	3.07	0.43
26:14:468:G:N7	54:L5:39:ARG:NH2	2.62	0.43
1:13:417:C:H2'	1:13:418:C:C6	2.54	0.43
48:F5:46:LEU:C	48:F5:47:GLN:HG2	2.38	0.43
1:13:491:G:H2'	1:13:492:G:O4'	2.19	0.43
29:29:170:LEU:HD11	29:29:185:LYS:O	2.19	0.43
33:61:29:TYR:C	33:61:32:PRO:HD2	2.39	0.43
26:1H:1292:U:H2'	26:1H:1293:C:O4'	2.19	0.43
6:5E:82:ARG:HB2	6:5E:85:VAL:HG23	2.00	0.43
38:98:32:GLY:HA2	38:98:116:LEU:HD12	2.01	0.43
1:1G:722:A:C8	1:1G:724:G:H1'	2.54	0.43
26:1H:800:A:OP1	58:1H:3653:HOH:O	2.21	0.43
30:31:68:LYS:O	30:31:69:HIS:HB2	2.18	0.43
20:BI:59:ALA:HA	20:BI:62:LEU:HD12	1.99	0.43
4:32:113:SER:O	4:32:117:ALA:N	2.51	0.43
27:1J:36:C:N3	27:1J:49:C:O2'	2.46	0.43
26:14:1782:C:H1'	26:14:2609:U:H5''	2.00	0.43
35:25:35:VAL:HG12	35:25:62:VAL:O	2.19	0.43
1:1G:1228:C:H5'	13:4A:114:ARG:HB3	2.01	0.43
1:1G:965:A:C2	1:1G:969:A:C2	3.07	0.43
1:1G:1246:C:H2'	1:1G:1247:U:O4'	2.18	0.43
55:Q8:47:LYS:HZ2	55:Q8:47:LYS:HA	1.83	0.43
18:9A:35:ARG:HE	18:9A:35:ARG:HB2	1.50	0.43
4:3E:154:ASN:ND2	4:3E:154:ASN:H	2.17	0.43
26:14:2582:G:H2'	26:14:2582:G:N3	2.34	0.43
2:1E:156:LYS:HA	2:1E:156:LYS:HD3	1.58	0.43
19:AA:12:ASP:HB3	19:AA:38:SER:HA	2.00	0.43
1:1G:977:A:OP1	14:5A:31:ARG:HD3	2.19	0.43
1:13:821:G:C2	1:13:880:C:N3	2.87	0.43
1:13:1530:G:H2'	1:13:1531:A:H8	1.84	0.43
1:1G:332:G:OP2	20:BA:10:LEU:HD12	2.19	0.43
36:78:19:VAL:CB	36:78:27:HIS:HB2	2.47	0.42
1:1G:974:A:P	14:5A:41:ARG:HH12	2.42	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:1142(A):A:C5	26:14:1144:G:C5	3.07	0.42
26:1H:120:U:C5	26:1H:149:A:N6	2.87	0.42
52:J5:16:ARG:HD2	52:J5:20:ARG:NH1	2.33	0.42
26:14:850:C:O3'	50:H5:49:LYS:HE2	2.19	0.42
55:M5:40:GLU:CA	55:M5:43:GLN:HB2	2.42	0.42
3:22:9:GLY:HA3	14:5A:49:HIS:HA	2.01	0.42
2:12:131:PRO:HG2	2:12:134:GLU:HB2	2.01	0.42
26:1H:2317:C:C2'	26:1H:2318:G:H5'	2.49	0.42
1:1G:468:A:H2'	1:1G:474:G:H5'	2.00	0.42
46:H8:128:VAL:HB	46:H8:161:VAL:HG21	2.01	0.42
38:98:79:LEU:HA	38:98:83:ILE:HG13	2.01	0.42
1:13:1124:G:H5'	10:1I:35:SER:HB2	2.01	0.42
55:Q8:9:GLY:HA2	55:Q8:12:LYS:HB2	2.00	0.42
26:14:747:U:P	52:J5:3:LYS:HD3	2.59	0.42
3:22:72:LYS:NZ	3:22:75:VAL:HG23	2.33	0.42
26:14:2527:C:C4	26:14:2528:U:C5	3.07	0.42
26:14:580:C:H2'	26:14:581:C:C6	2.54	0.42
46:D5:91:LEU:HD22	46:D5:96:VAL:HG21	2.00	0.42
8:72:12:ARG:NH2	8:72:27:PRO:HD3	2.34	0.42
26:1H:2689:U:H4'	26:1H:2690:C:H5'	2.01	0.42
36:35:78:PRO:HB3	36:35:111:ARG:HH21	1.83	0.42
27:1J:9:G:OP1	39:65:25:ARG:NH2	2.52	0.42
13:4I:82:MET:C	13:4I:84:ILE:H	2.23	0.42
26:1H:1009:A:OP2	34:58:37:LYS:NZ	2.51	0.42
6:5E:55:ASP:HB2	6:5E:86:ARG:HH12	1.83	0.42
1:1G:571:U:O2	1:1G:918:A:H5'	2.19	0.42
2:12:142:LEU:HD23	2:12:142:LEU:O	2.19	0.42
26:14:2467:C:H2'	26:14:2468:G:O4'	2.18	0.42
30:39:181:LEU:HD21	30:39:186:ILE:HD11	2.01	0.42
26:14:2862:G:H2'	26:14:2863:C:C6	2.54	0.42
38:98:29:LEU:HB3	38:98:75:LEU:HD11	2.01	0.42
16:7A:15:PRO:C	16:7A:16:HIS:HD2	2.22	0.42
1:13:490:G:O2'	1:13:491:G:H5'	2.18	0.42
26:14:1921:G:H2'	26:14:1922:G:H8	1.84	0.42
29:21:96:PHE:O	29:21:175:VAL:HG11	2.18	0.42
26:14:1312:U:H4'	26:14:1313:U:O5'	2.19	0.42
1:13:883:C:O2'	1:13:884:U:H5'	2.19	0.42
26:14:979:G:H3'	26:14:980:A:C5'	2.49	0.42
36:78:132:LYS:HB3	36:78:132:LYS:HE3	1.83	0.42
20:BI:29:LYS:HB2	20:BI:29:LYS:HE3	1.81	0.42
44:F8:2:LYS:O	44:F8:2:LYS:HE2	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:402:A:O5'	26:14:402:A:H8	2.02	0.42
26:14:1813:G:H1'	28:19:50:THR:OG1	2.18	0.42
2:12:50:GLU:HB3	2:12:200:ILE:O	2.19	0.42
26:1H:2393:A:C2'	26:1H:2394:C:H5'	2.49	0.42
43:A5:88:ARG:HB3	43:A5:92:ARG:HB3	2.00	0.42
26:1H:1556:C:H2'	26:1H:1557:C:C6	2.53	0.42
55:Q8:53:PRO:CB	55:Q8:56:GLU:H	2.32	0.42
26:1H:456:C:C4	44:F8:69:TYR:CE1	3.07	0.42
26:1H:57:C:H2'	26:1H:58:G:O4'	2.19	0.42
1:13:738:C:H2'	1:13:739:C:C6	2.53	0.42
26:14:631:A:N3	26:14:2415:G:O2'	2.34	0.42
26:14:1754:C:OP2	40:75:113:LYS:NZ	2.52	0.42
26:1H:141:A:OP2	26:1H:141(A):C:N4	2.33	0.42
1:13:1182:G:H4'	1:13:1183:A:C5'	2.49	0.42
4:32:18:LYS:HD2	4:32:20:TYR:CE1	2.55	0.42
27:1J:42:C:N4	27:1J:43:C:C4	2.87	0.42
28:11:69:ARG:NH2	28:11:128:GLY:O	2.39	0.42
27:1J:104:A:O4'	46:D5:29:TYR:HE2	2.02	0.42
42:95:80:GLN:HG3	42:95:81:TYR:N	2.25	0.42
30:31:9:ILE:HG12	30:31:10:PRO:HD2	2.01	0.42
1:1G:1321:C:H4'	13:4A:87:TYR:CZ	2.53	0.42
45:G8:74:PRO:O	45:G8:82:PRO:HD2	2.18	0.42
13:4A:39:ILE:HG12	13:4A:55:ARG:HH21	1.85	0.42
26:14:1945:G:C4	26:14:1946:U:C5	3.07	0.42
26:14:2360:A:H2'	26:14:2361:A:O4'	2.19	0.42
4:32:15:GLU:OE1	4:32:59:ARG:NH2	2.34	0.42
26:1H:569:U:O4	26:1H:570:G:C6	2.72	0.42
26:1H:270(E):G:N2	26:1H:270(U):C:N3	2.59	0.42
26:14:746:A:H2'	26:14:2612:C:H5''	2.00	0.42
26:1H:1591:G:O2'	26:1H:1592:C:H5'	2.19	0.42
26:1H:2148:G:H2'	26:1H:2149:G:O4'	2.19	0.42
34:58:46:VAL:CG1	34:58:48:MET:HG3	2.50	0.42
22:1K:70:G:H4'	22:1K:70:G:OP1	2.18	0.42
1:13:474:G:H5''	16:7I:81:ARG:NE	2.35	0.42
1:13:474:G:C6	1:13:475:G:C6	3.06	0.42
1:13:1312:G:H5'	19:AI:6:LYS:HD2	2.01	0.42
45:G8:45:VAL:N	45:G8:63:LYS:O	2.47	0.42
26:1H:931:G:H4'	50:L8:24:LYS:NZ	2.34	0.42
26:1H:1766:U:H2'	26:1H:1767:C:C6	2.54	0.42
34:58:76:SER:O	34:58:78:TYR:N	2.51	0.42
1:1G:1292:U:H2'	1:1G:1293:G:C8	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
40:75:18:ASP:N	40:75:18:ASP:OD1	2.50	0.42
26:1H:2748:A:H2	32:51:63:SER:HB3	1.82	0.42
26:14:1805:U:H2'	26:14:1806:C:H6	1.84	0.42
24:3K:34:G:H2'	24:3K:35:A:C8	2.54	0.42
28:11:119:ALA:HB1	28:11:130:ALA:HB3	2.00	0.42
13:4I:58:GLU:O	13:4I:62:ASN:HB2	2.19	0.42
1:13:380:G:N2	1:13:384:G:C6	2.87	0.42
1:13:1350:A:C5	1:13:1351:U:C4	3.08	0.42
1:13:933:G:OP2	7:6E:3:ARG:HB2	2.19	0.42
16:7I:83:GLU:HB3	16:7I:84:ALA:H	1.54	0.42
26:1H:381:G:C4	26:1H:394:A:C2	3.07	0.42
2:1E:82:ARG:HH21	2:1E:150:SER:HB3	1.84	0.42
26:14:2097:C:H2'	26:14:2098:U:C6	2.54	0.42
1:1G:1511:G:H8	1:1G:1511:G:O5'	2.02	0.42
22:1K:62:C:H2'	22:1K:63:G:C8	2.53	0.42
35:25:8:LEU:HD13	35:25:82:ASN:CB	2.49	0.42
1:13:645:C:H2'	1:13:646:U:O4'	2.18	0.42
26:1H:1344:G:C2	26:1H:1385:G:C8	3.08	0.42
26:14:631:A:H2'	26:14:632:A:O4'	2.18	0.42
26:1H:191:A:H2'	26:1H:192:C:C6	2.54	0.42
4:32:18:LYS:HB2	4:32:31:CYS:SG	2.58	0.42
48:J8:91:LYS:O	48:J8:93:GLU:N	2.52	0.42
29:29:16:ARG:HH11	29:29:16:ARG:HG3	1.85	0.42
45:G8:76:CYS:HB2	45:G8:97:ARG:HG2	2.00	0.42
1:1G:555:C:H2'	1:1G:556:C:C6	2.54	0.42
35:68:78:ARG:HH21	40:B8:103:ARG:HH21	1.68	0.42
1:13:1145:C:H4'	1:13:1146:A:C8	2.46	0.42
1:1G:999:U:H2'	1:1G:1000:A:H8	1.82	0.42
29:29:57:LYS:HD3	29:29:57:LYS:HA	1.61	0.42
26:1H:1341:U:H4'	26:1H:1342:A:OP2	2.18	0.42
5:42:5:ASP:HA	5:42:63:ARG:HH12	1.83	0.42
5:42:6:PHE:HD2	5:42:63:ARG:HD3	1.83	0.42
3:22:47:LEU:HB3	3:22:52:LEU:HD13	2.01	0.42
9:8E:40:LEU:HD11	9:8E:70:LYS:HD3	2.01	0.42
1:1G:987:G:O5'	1:1G:987:G:H8	2.02	0.42
5:4E:147:ASP:HA	5:4E:150:ARG:NH2	2.35	0.42
1:1G:160:A:H1'	1:1G:344:A:N7	2.34	0.42
1:13:555:C:H2'	1:13:556:C:C6	2.54	0.42
36:78:147:LEU:HD12	36:78:147:LEU:HA	1.81	0.42
49:K8:24:LEU:HA	49:K8:24:LEU:HD23	1.84	0.42
8:7E:16:ALA:CB	8:7E:24:THR:HG21	2.48	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
29:21:152:LYS:HG2	34:58:78:TYR:CE1	2.54	0.42
26:1H:1542:G:OP2	26:1H:1543:A:O2'	2.30	0.42
26:1H:16:G:H2'	26:1H:17:G:H8	1.84	0.42
1:13:321:A:H62	1:13:328:C:H1'	1.84	0.42
42:D8:21:ARG:HG2	42:D8:91:TYR:HE2	1.82	0.42
1:1G:1134:G:H2'	1:1G:1135:U:O4'	2.19	0.42
26:1H:902:C:H2'	26:1H:903:C:C6	2.54	0.42
4:32:71:SER:OG	4:32:74:GLN:HG3	2.20	0.42
26:14:948:G:C2	26:14:970:C:O2	2.73	0.42
30:39:64:ILE:HG13	30:39:65:TRP:CD1	2.54	0.42
7:6E:66:VAL:O	7:6E:70:LYS:HG3	2.19	0.42
2:12:109:SER:HA	2:12:112:VAL:HG23	2.02	0.42
26:14:387:U:H4'	26:14:388:G:O5'	2.19	0.42
38:98:84:ALA:HB3	38:98:85:PRO:HD3	2.00	0.42
30:31:116:ASP:O	30:31:120:GLU:HG3	2.18	0.42
54:L5:26:GLY:O	54:L5:30:VAL:HG23	2.19	0.42
3:2E:137:ALA:O	3:2E:141:VAL:HG23	2.19	0.42
32:51:88:LEU:O	32:51:88:LEU:HD12	2.17	0.42
26:1H:339:U:H6	26:1H:339:U:O5'	2.02	0.42
19:AI:30:LEU:HD13	19:AI:30:LEU:H	1.85	0.42
50:H5:5:LYS:HB3	50:H5:5:LYS:HE3	1.60	0.42
37:45:58:PHE:CD2	37:45:61:GLY:HA3	2.54	0.42
1:13:734:G:C2	1:13:735:C:C2	3.08	0.42
2:1E:185:ILE:HA	2:1E:199:TYR:O	2.19	0.42
1:13:1391:U:H2'	1:13:1392:G:H8	1.76	0.42
30:39:185:ASP:CG	30:39:188:ARG:HH21	2.21	0.42
1:1G:974:A:H5'	1:1G:975:A:OP1	2.18	0.42
5:42:107:ARG:O	5:42:110:LEU:N	2.53	0.42
1:1G:680:C:N3	1:1G:710:G:N2	2.53	0.42
26:1H:142:G:H1'	44:F8:37:THR:CG2	2.41	0.42
26:14:2312:U:C5	26:14:2313:C:C4	3.08	0.42
39:A8:89:ARG:O	39:A8:89:ARG:CG	2.66	0.42
1:1G:673:G:O3'	6:52:87:ARG:NH2	2.52	0.42
24:3K:5:G:H2'	24:3K:6:G:O4'	2.19	0.42
28:11:206:LEU:O	28:11:211:ARG:HD3	2.20	0.42
26:14:2331:G:O2'	47:E5:43:THR:HB	2.19	0.42
12:3I:53:ARG:HH12	12:3I:92:ASP:CB	2.32	0.42
31:41:99:MET:HG3	31:41:100:TRP:N	2.34	0.42
1:13:524:G:H2'	1:13:525:C:C5	2.53	0.42
1:1G:1117:G:O3'	9:82:104:ARG:HD2	2.19	0.42
26:14:249:C:H4'	26:14:250:G:O5'	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:1H:286:C:H2'	26:1H:287:C:H6	1.84	0.42
41:85:91:ASP:OD2	41:85:96:ALA:HB2	2.19	0.42
1:13:427:U:H3'	1:13:428:G:H2'	2.01	0.42
1:13:324:G:H22	1:13:327:A:P	2.41	0.42
33:69:75:LEU:HA	33:69:139:GLN:OE1	2.19	0.42
26:1H:1431:U:C2	26:1H:1563:G:N2	2.87	0.42
13:4A:31:LYS:HA	13:4A:34:LEU:HD12	2.01	0.42
12:3I:11:VAL:HG13	17:8I:29:HIS:ND1	2.34	0.42
13:4I:27:LYS:HD3	13:4I:27:LYS:HA	1.71	0.42
26:14:871:U:OP1	37:45:5:ARG:HG2	2.18	0.42
28:19:12:SER:HB2	28:19:208:LYS:HB3	2.00	0.42
26:1H:763:G:O2'	26:1H:764:A:H3'	2.19	0.42
49:K8:3:LEU:O	49:K8:6:VAL:HG22	2.19	0.42
1:13:222:U:C2	1:13:223:U:C5	3.07	0.42
1:13:313:A:H2'	1:13:314:C:C6	2.54	0.42
40:75:45:PHE:CE1	40:75:65:LYS:HG2	2.55	0.42
1:1G:485:G:O2'	1:1G:486:U:O5'	2.36	0.42
26:1H:363(B):G:H2'	26:1H:363(C):G:C8	2.52	0.42
55:Q8:36:LYS:O	55:Q8:40:GLU:HB2	2.19	0.42
37:88:65:PHE:HD1	37:88:105:GLU:O	2.02	0.42
6:52:96:PRO:HB3	18:9A:30:ASP:OD2	2.20	0.42
5:42:80:ILE:HG13	8:72:104:ARG:HH21	1.84	0.42
1:13:1277:C:O2'	1:13:1279:A:H1'	2.19	0.42
26:14:1338:G:N3	26:14:1393:A:H2	2.17	0.42
26:14:1654:A:C1'	26:14:2823:A:H5'	2.49	0.42
1:1G:967:C:H2'	1:1G:968:A:C8	2.54	0.42
34:15:28:THR:HG22	34:15:29:LYS:N	2.34	0.42
30:39:82:ILE:HG13	30:39:82:ILE:H	1.50	0.42
1:1G:5:U:O2'	4:32:84:LYS:HG3	2.19	0.42
26:1H:557:U:C2	26:1H:558:G:C8	3.08	0.42
4:32:93:PHE:CZ	4:32:97:LEU:HD11	2.54	0.42
10:1I:31:GLY:HA3	10:1I:81:THR:CG2	2.48	0.42
26:1H:216:A:C4	26:1H:432:A:C2	3.07	0.42
26:14:1812:A:H2'	26:14:1813:G:C8	2.54	0.42
26:1H:343:C:O2'	26:1H:344:G:H5'	2.19	0.42
26:14:2852:G:H2'	26:14:2853:C:O4'	2.20	0.42
26:14:2579:C:H2'	26:14:2580:U:O4'	2.19	0.42
26:14:1465:G:C4	26:14:1466:G:C8	3.07	0.42
39:65:7:TYR:CZ	39:65:91:PRO:HG3	2.54	0.42
3:22:65:ALA:HA	3:22:100:ALA:HB3	2.01	0.42
26:1H:1812:A:O2'	28:11:45:ASN:N	2.53	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
30:39:24:LEU:HD11	30:39:119:ARG:HB2	2.01	0.42
26:1H:271(B):G:H4'	26:1H:271(C):U:C5'	2.49	0.42
1:13:669:U:C2	1:13:670:G:C8	3.07	0.42
29:29:97:LYS:O	29:29:100:GLU:HG3	2.20	0.42
1:13:406:G:H5'	4:3E:5:ILE:HD13	2.00	0.42
1:1G:44:G:N2	1:1G:398:C:O2	2.40	0.42
26:14:1950:G:C2	26:14:1951:U:C5	3.06	0.42
46:H8:103:ARG:HG3	46:H8:136:PHE:CD2	2.54	0.42
32:59:105:LEU:HG	32:59:113:VAL:HG13	2.01	0.42
26:1H:1034:G:C5	26:1H:1035:U:C5	3.07	0.42
42:D8:66:ARG:HH21	42:D8:88:ARG:HH11	1.65	0.42
17:8I:52:LYS:HD2	17:8I:55:ASP:OD1	2.19	0.42
26:14:1051:G:H8	26:14:1051:G:OP2	2.03	0.42
26:14:2003:G:C6	26:14:2004:G:N7	2.88	0.42
26:1H:900:A:H5'	26:1H:901:A:P	2.59	0.42
36:78:60:MET:HA	55:Q8:13:ARG:NH1	2.34	0.42
1:1G:21:G:H2'	1:1G:22:G:C8	2.54	0.42
20:BI:73:HIS:HB3	20:BI:74:LYS:CD	2.44	0.42
27:1J:56:G:H4'	27:1J:57:A:N7	2.35	0.42
30:39:117:ARG:NH1	36:35:1:MET:H2	2.08	0.42
10:1I:54:PHE:CG	10:1I:55:LYS:HG2	2.54	0.42
26:1H:996:A:O3'	41:C8:92:ARG:HG2	2.18	0.42
50:L8:7:LYS:O	50:L8:54:VAL:HA	2.20	0.42
27:16:40:U:C5	51:M8:2:LYS:HG2	2.55	0.42
30:39:122:LYS:HD2	30:39:191:ARG:HE	1.84	0.42
1:1G:984:C:H2'	1:1G:985:C:H6	1.84	0.42
2:1E:53:ARG:NH1	2:1E:200:ILE:HD12	2.29	0.42
1:1G:428:G:C5	1:1G:430:A:C6	3.08	0.42
26:1H:2684:U:O2'	35:68:68:GLU:HG3	2.18	0.42
26:14:971:C:H2'	26:14:972:G:O4'	2.19	0.42
3:2E:177:THR:HB	3:2E:180:ALA:HB2	2.01	0.42
26:1H:2859:G:C6	26:1H:2860:A:N6	2.87	0.42
9:8E:46:ALA:HB2	9:8E:74:ILE:HG23	2.00	0.42
55:Q8:33:ASN:HB3	55:Q8:34:TRP:H	1.60	0.42
26:1H:1590:U:C2	26:1H:1591:G:N7	2.87	0.42
1:13:1052:U:O2'	1:13:1055:A:OP2	2.32	0.42
28:11:46:GLN:H	28:11:46:GLN:HG3	1.44	0.42
30:39:205:ARG:HB2	30:39:205:ARG:NH1	2.33	0.42
26:14:1386:C:OP2	26:14:1396:U:C5	2.73	0.42
20:BI:35:THR:O	20:BI:38:LYS:HB2	2.19	0.42
1:1G:91:C:H2'	1:1G:92:G:C8	2.55	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:13:688:G:H2'	1:13:689:C:C6	2.52	0.42
29:29:116:VAL:HG21	29:29:122:PHE:CE2	2.55	0.42
1:1G:865:A:H2	1:1G:918:A:H4'	1.85	0.42
26:1H:2335:A:C8	26:1H:2337:G:C6	3.07	0.42
20:BA:67:ALA:HA	20:BA:73:HIS:H	1.83	0.42
41:C8:49:HIS:HA	41:C8:52:ARG:HG2	2.01	0.42
1:1G:56:U:H2'	1:1G:57:G:H8	1.84	0.42
1:13:1117:G:H5''	9:8E:104:ARG:NH1	2.33	0.42
1:1G:4:U:C4	8:72:102:ARG:HD2	2.54	0.42
26:1H:1952:A:H5''	26:1H:1953:A:OP2	2.19	0.42
26:14:2022:U:OP2	52:J5:15:ARG:NH2	2.52	0.42
53:K5:11:LEU:HD13	53:K5:26:ASN:HD22	1.84	0.42
5:42:31:LEU:HD23	5:42:45:PHE:HB2	2.02	0.42
26:1H:2677:G:N2	26:1H:2731:G:C4	2.87	0.42
26:14:2506:U:H4'	26:14:2507:C:OP1	2.15	0.42
15:6A:45:VAL:HB	15:6A:46:HIS:CD2	2.55	0.42
33:69:29:TYR:C	33:69:32:PRO:HD2	2.40	0.42
26:1H:2516:G:O2'	26:1H:2517:C:H5'	2.19	0.42
1:1G:1197:G:H2'	1:1G:1197:G:N3	2.35	0.42
45:G8:88:LYS:HA	45:G8:88:LYS:HD3	1.53	0.42
16:7A:43:LYS:HE2	16:7A:43:LYS:HB2	1.66	0.42
34:15:127:ASP:N	34:15:127:ASP:OD1	2.53	0.42
9:8E:92:TYR:HD1	9:8E:92:TYR:HA	1.69	0.42
1:1G:986:A:H1'	19:AA:54:GLY:O	2.18	0.42
13:4I:32:GLU:OE2	13:4I:33:ALA:N	2.52	0.42
26:1H:2364:C:H4'	47:I8:56:ASP:OD1	2.20	0.42
26:1H:2078:C:H2'	26:1H:2079:U:C6	2.54	0.42
26:1H:512:G:C8	58:1H:4190:HOH:O	2.57	0.42
1:13:452:A:O2'	1:13:453:A:O4'	2.34	0.42
26:1H:442:G:C6	26:1H:444:C:N4	2.87	0.42
26:14:2129:C:H2'	26:14:2130:U:O4'	2.19	0.42
30:31:9:ILE:HD12	30:31:125:LEU:HG	2.01	0.42
32:51:4:ILE:CD1	32:51:4:ILE:H	2.26	0.42
4:32:8:VAL:O	4:32:11:LEU:N	2.39	0.42
26:14:2262:U:H4'	26:14:2328:A:C2	2.54	0.42
31:41:131:TYR:HB3	31:41:159:VAL:CG2	2.50	0.42
17:8I:22:LEU:HD13	17:8I:41:LYS:HG3	2.01	0.42
26:1H:242:G:H5'	55:Q8:61:LEU:HD22	2.02	0.42
1:1G:1172:C:H2'	1:1G:1173:G:C8	2.55	0.42
13:4I:11:ARG:HD3	13:4I:46:LYS:NZ	2.33	0.42
26:14:1639:U:H2'	26:14:1640:C:H5'	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
38:98:63:ARG:HG2	38:98:67:LEU:HD23	2.01	0.42
29:21:13:ARG:HH11	29:21:13:ARG:CG	2.33	0.42
1:13:412:A:N6	4:3E:35:ARG:HA	2.34	0.42
1:1G:246:A:N6	1:1G:281:G:H1'	2.35	0.42
1:13:1286:A:H5''	21:1F:26:LYS:CG	2.48	0.42
33:69:141:LYS:HE2	33:69:141:LYS:HB2	1.88	0.42
26:14:2014:A:H2'	26:14:2015:A:C8	2.55	0.42
26:14:654(C):G:C2	26:14:654(S):G:C2	3.07	0.42
53:K5:9:LEU:N	53:K5:27:LYS:HG3	2.34	0.42
26:1H:1590:U:N3	26:1H:1591:G:N7	2.68	0.42
26:14:854:G:H2'	26:14:855:G:C8	2.54	0.42
26:1H:993:G:OP1	41:C8:50:ARG:NH2	2.52	0.42
1:13:179:A:H2'	1:13:180:U:C6	2.53	0.42
4:3E:70:ILE:HG23	4:3E:75:PHE:HB2	1.99	0.42
26:14:235:U:H2'	26:14:236:C:C6	2.54	0.42
28:11:61:LEU:HA	28:11:61:LEU:HD13	1.72	0.42
18:9I:70:ILE:HG23	18:9I:79:LEU:HD12	2.02	0.42
27:1J:83:G:H4'	50:H5:52:HIS:CG	2.55	0.42
2:12:142:LEU:HA	2:12:145:LEU:HB2	2.01	0.42
29:29:37:ARG:HD3	29:29:44:TYR:CZ	2.54	0.42
37:45:110:THR:OG1	37:45:112:GLU:OE2	2.29	0.42
19:AA:29:ARG:HH11	19:AA:48:THR:HG23	1.84	0.42
26:1H:319:C:H2'	26:1H:320:A:C8	2.54	0.42
39:65:106:ARG:H	39:65:106:ARG:HG3	1.48	0.42
26:14:754:C:H2'	26:14:755:C:H6	1.84	0.42
32:59:103:LEU:HD22	32:59:123:PHE:CE1	2.54	0.42
27:16:89(A):A:C5	27:16:90:C:H1'	2.54	0.42
5:4E:81:GLU:HG2	5:4E:90:VAL:HG23	2.02	0.42
45:G8:43:ASN:ND2	45:G8:67:LEU:HD11	2.35	0.42
26:14:2031:A:N3	26:14:2455:G:O2'	2.45	0.42
26:14:2212:A:H1'	26:14:2215:G:C5	2.54	0.42
29:21:134:ILE:HD12	29:21:134:ILE:C	2.39	0.42
1:1G:1072:G:C2	1:1G:1073:U:C2	3.07	0.42
20:BA:10:LEU:HD13	20:BA:10:LEU:O	2.20	0.42
37:45:58:PHE:HZ	37:45:106:VAL:HG11	1.85	0.42
26:14:1499:C:H2'	26:14:1500:G:H8	1.84	0.42
1:1G:637:G:C2	1:1G:638:G:C4	3.08	0.42
22:1K:54:5MU:H2'	22:1K:55:PSU:O4'	2.19	0.42
26:14:1253:A:OP1	58:14:3962:HOH:O	2.21	0.42
35:68:24:VAL:HB	35:68:33:ALA:HB2	2.01	0.42
1:13:949:A:C2	1:13:1233:G:N3	2.88	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:7E:75:ARG:HE	8:7E:75:ARG:HB2	1.50	0.42
26:14:205:G:O2'	26:14:206:U:OP2	2.36	0.42
26:14:1682:G:H2'	26:14:1683:C:C6	2.55	0.42
26:14:2436:G:C5	26:14:2437:U:C5	3.08	0.42
1:13:8:A:H5''	5:4E:120:THR:O	2.19	0.42
26:1H:496:G:H2'	26:1H:497:A:H5'	2.00	0.42
28:11:27:THR:OG1	28:11:27:THR:O	2.34	0.42
18:9I:59:SER:OG	18:9I:60:ALA:N	2.51	0.42
33:61:75:LEU:HB3	33:61:105:HIS:HD2	1.85	0.42
26:1H:660:G:O3'	30:31:38:ARG:NH2	2.52	0.42
26:1H:2155:G:H2'	26:1H:2156:G:H5'	2.00	0.42
1:1G:1281:U:H3'	1:1G:1282:C:H5	1.81	0.42
9:82:9:ARG:HG2	9:82:14:VAL:HG22	2.01	0.42
55:Q8:7:HIS:CB	55:Q8:58:ILE:HG23	2.45	0.42
50:L8:35:ARG:HB3	50:L8:37:LEU:HD22	2.02	0.42
7:62:114:ARG:H	7:62:114:ARG:HG2	1.51	0.42
26:14:38:A:H1'	30:39:48:THR:HB	2.02	0.42
27:1J:43:C:H4'	31:49:98:ARG:HH12	1.84	0.42
27:16:42:C:O2'	31:41:67:LYS:O	2.29	0.42
26:14:619:G:OP2	26:14:620:G:N2	2.40	0.42
16:7A:81:ARG:HD2	16:7A:83:GLU:OE2	2.19	0.42
1:13:127:G:H4'	17:8I:2:PRO:HD2	2.01	0.42
26:1H:1324:G:C5	26:1H:1328:G:O6	2.73	0.42
26:1H:657:U:H2'	26:1H:658:C:C6	2.54	0.42
28:11:136:ILE:HG12	28:11:136:ILE:H	1.61	0.42
29:29:11:MET:HA	29:29:24:THR:HA	2.01	0.42
1:1G:939:G:H1	1:1G:1344:C:H42	1.67	0.42
5:4E:37:ARG:HA	5:4E:114:GLY:H	1.84	0.42
3:2E:6:HIS:HA	3:2E:7:PRO:HD2	1.88	0.42
26:1H:1693:U:H4'	26:1H:1694:C:OP2	2.20	0.42
4:3E:96:LEU:HD12	4:3E:139:ARG:HD3	2.01	0.42
30:31:164:ARG:HG3	30:31:175:THR:OG1	2.19	0.42
26:1H:2481:G:HO2'	26:1H:2482:G:P	2.42	0.42
36:35:11:GLY:HA2	36:35:13:ASN:OD1	2.19	0.42
1:1G:746:A:H2'	1:1G:747:C:C6	2.54	0.42
50:L8:26:LEU:CB	50:L8:28:LEU:HD12	2.49	0.42
33:61:81:VAL:HG23	33:61:143:SER:O	2.20	0.42
17:8I:74:LEU:HB3	17:8I:75:ARG:HG2	2.00	0.42
26:14:2642:G:P	34:15:76:SER:HG	2.42	0.42
26:1H:1444:G:N2	26:1H:1548:C:C2	2.88	0.42
1:13:784:C:H2'	1:13:785:G:C8	2.55	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
33:61:128:LEU:O	33:61:138:ILE:N	2.52	0.42
26:14:1644:C:O2'	26:14:1645:G:H5'	2.19	0.42
1:13:373:A:C2	1:13:374:A:C8	3.07	0.42
1:1G:1442:G:H2'	1:1G:1443:G:H5''	2.01	0.42
49:K8:8:LYS:NZ	49:K8:12:GLU:HG2	2.34	0.42
26:14:61:G:O5'	26:14:61:G:H8	2.02	0.42
49:G5:16:LEU:HG	49:G5:16:LEU:O	2.19	0.42
46:D5:52:SER:C	46:D5:54:HIS:H	2.21	0.42
53:K5:52:VAL:HG22	53:K5:53:LYS:H	1.85	0.42
26:14:1509:C:H5'	26:14:1510:A:O4'	2.19	0.42
39:65:47:THR:HG22	39:65:48:LEU:H	1.84	0.42
30:31:184:TYR:HE1	36:78:3:LEU:HD11	1.85	0.42
26:14:2459:A:C4	26:14:2460:U:C5	3.08	0.42
26:14:2854:G:C2	26:14:2864:G:C2	3.07	0.42
1:13:1386:G:O2'	1:13:1387:G:H5'	2.20	0.42
26:14:2865:U:C4	26:14:2866:U:C4	3.07	0.42
26:1H:470:A:C2	26:1H:471:A:C4	3.07	0.42
26:1H:631:A:H61	26:1H:2402:C:N4	2.18	0.42
26:14:805:G:H4'	36:35:38:GLN:HB2	2.01	0.42
26:14:1403:C:OP1	26:14:1522:G:N2	2.49	0.42
26:14:2168:G:N3	26:14:2168:G:H2'	2.34	0.42
51:I5:14:ILE:HB	51:I5:24:THR:HG21	2.01	0.42
46:D5:74:VAL:HA	46:D5:86:VAL:HG22	2.02	0.42
55:M5:8:LYS:HD3	55:M5:8:LYS:HA	1.62	0.42
12:3A:123:LYS:HG2	12:3A:123:LYS:H	1.44	0.42
18:9A:84:LYS:HG2	18:9A:84:LYS:H	1.50	0.42
2:1E:74:LYS:HD2	2:1E:74:LYS:H	1.85	0.42
26:1H:139:G:N3	26:1H:141:A:N1	2.67	0.42
27:1J:40:U:H1'	27:1J:46:A:N1	2.34	0.42
14:5I:29:ARG:HH21	14:5I:41:ARG:NH1	2.18	0.42
1:1G:1048:G:N2	1:1G:1210:C:N3	2.67	0.42
1:1G:1206:G:C6	1:1G:1207:G:C5	3.07	0.42
26:14:2689:U:H5''	26:14:2713:A:C2	2.54	0.42
1:13:1005:A:H5''	1:13:1038:C:H1'	2.01	0.42
45:C5:101:LYS:C	45:C5:102:CYS:SG	2.98	0.42
40:B8:9:LEU:HD23	40:B8:9:LEU:HA	1.90	0.42
30:31:9:ILE:HG22	30:31:20:LEU:O	2.19	0.42
30:39:122:LYS:CB	30:39:191:ARG:HB2	2.49	0.42
43:E8:12:ILE:HD13	43:E8:17:VAL:HB	2.02	0.42
40:B8:76:PHE:HB3	40:B8:83:ILE:HD11	2.02	0.42
26:14:2244:U:O5'	26:14:2244:U:H6	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
46:D5:30:ASN:HB3	46:D5:90:VAL:HB	2.02	0.42
1:13:232:G:H1'	1:13:262:A:N1	2.34	0.42
1:13:232:G:H2'	1:13:233:C:C6	2.55	0.42
53:O8:20:ASN:C	53:O8:21:TYR:CG	2.92	0.42
1:13:648:A:N6	1:13:649:G:O6	2.53	0.42
30:31:114:VAL:HG21	30:31:202:PHE:CZ	2.54	0.42
1:1G:691:G:O6	11:2A:55:LYS:NZ	2.39	0.42
35:25:10:VAL:HG23	35:25:10:VAL:O	2.20	0.42
18:9A:22:VAL:CG1	18:9A:56:THR:HA	2.49	0.42
3:22:37:GLN:O	3:22:41:GLY:N	2.50	0.42
43:E8:86:LEU:HD12	43:E8:87:PRO:CD	2.50	0.42
28:19:37:LEU:HA	28:19:38:LYS:CB	2.50	0.42
1:13:577:G:O2'	1:13:816:A:H2'	2.20	0.42
26:1H:250:G:C6	26:1H:251:A:C6	3.08	0.42
4:3E:134:ASP:HB2	4:3E:135:LEU:HD13	2.01	0.42
3:22:156:ARG:HB2	3:22:159:GLY:HA2	2.02	0.42
26:1H:270(N):G:OP2	33:61:57:ARG:NH2	2.52	0.42
6:52:71:ARG:H	6:52:71:ARG:HG2	1.51	0.42
32:59:135:GLY:HA3	32:59:141:VAL:HG22	2.02	0.42
26:1H:2309:A:C4	26:1H:2310:A:H8	2.37	0.42
1:13:1308:U:OP1	13:4I:98:VAL:HG12	2.20	0.42
18:9A:19:LYS:HA	18:9A:19:LYS:HD2	1.94	0.42
32:51:149:ARG:HG3	32:51:149:ARG:HH11	1.85	0.42
28:19:106:ILE:O	28:19:108:PRO:HD3	2.19	0.42
42:D8:66:ARG:NH2	42:D8:88:ARG:HH11	2.17	0.42
11:2I:91:ARG:O	11:2I:94:ALA:HB3	2.19	0.42
26:1H:827:U:H5'	26:1H:828:U:O5'	2.19	0.42
11:2I:18:ARG:HB3	11:2I:33:THR:OG1	2.19	0.42
42:D8:14:VAL:HA	42:D8:18:LEU:HD12	2.02	0.42
9:82:84:ALA:O	9:82:87:GLN:HB3	2.20	0.42
39:65:24:LEU:HB2	39:65:85:VAL:HG12	2.02	0.42
48:F5:21:ARG:HD3	48:F5:35:THR:HG21	2.02	0.42
10:1I:32:ALA:HB1	10:1I:76:ASN:OD1	2.20	0.42
26:14:1133:U:O2	26:14:1137:G:H5'	2.19	0.42
1:1G:131:C:H2'	1:1G:132:C:C6	2.55	0.42
18:9I:36:ASN:ND2	18:9I:39:VAL:HG21	2.34	0.42
35:68:90:GLN:O	35:68:91:LEU:HB2	2.18	0.42
26:1H:1414:G:C6	26:1H:1415:U:C4	3.07	0.42
5:4E:87:SER:HB3	5:4E:125:SER:O	2.20	0.42
26:14:452:G:N3	26:14:457:A:H2	2.16	0.42
26:1H:1006:C:C2	26:1H:1138:G:N2	2.87	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:1H:2577:A:C5'	26:1H:2578:G:H5'	2.46	0.42
1:1G:1149:C:OP1	9:82:9:ARG:HD3	2.20	0.42
26:14:1141:U:P	34:15:25:ARG:HH21	2.42	0.42
1:13:537:G:H2'	1:13:538:G:C8	2.55	0.42
26:1H:2052:G:C2	26:1H:2053:G:C8	3.08	0.42
26:1H:2128:C:H2'	26:1H:2129:C:C6	2.54	0.42
1:13:165:C:H2'	1:13:166:G:O4'	2.19	0.42
26:14:2074:U:P	58:14:3509:HOH:O	2.75	0.42
26:1H:2608:G:O5'	26:1H:2608:G:H8	2.03	0.42
27:1J:42:C:N3	31:49:91:ARG:NH2	2.67	0.42
26:14:2135:A:O2'	26:14:2160:G:H4'	2.19	0.42
19:AI:42:PRO:O	19:AI:45:VAL:HG22	2.20	0.42
26:14:2712:U:O2'	26:14:2712(A):A:P	2.77	0.42
26:14:275:G:H2'	26:14:276:A:C8	2.54	0.42
16:7I:4:ILE:HA	16:7I:20:VAL:O	2.19	0.42
1:13:266:G:N2	1:13:269:C:C5	2.88	0.42
1:13:101:A:C8	1:13:101:A:OP2	2.73	0.42
26:14:2361:A:N7	58:14:3994:HOH:O	2.37	0.42
13:4A:8:GLU:OE2	13:4A:10:PRO:HG3	2.20	0.42
2:1E:183:PRO:HA	2:1E:198:ASP:OD2	2.20	0.42
14:5I:9:LYS:HA	14:5I:12:ARG:HG2	2.02	0.42
1:13:375:U:O3'	16:7I:6:LEU:HB2	2.20	0.42
26:1H:537:C:H2'	26:1H:539:G:C8	2.55	0.42
42:D8:9:GLY:O	42:D8:10:LYS:HG3	2.20	0.42
35:25:47:ILE:HD12	35:25:47:ILE:HA	1.76	0.42
13:4A:78:ILE:HD13	13:4A:92:HIS:HE1	1.84	0.42
12:3A:84:LEU:C	12:3A:85:ILE:HD12	2.40	0.42
29:29:44:TYR:HE1	29:29:80:GLU:OE1	2.03	0.42
27:16:61:G:C6	27:16:62:C:C4	3.08	0.42
26:14:2079:U:H2'	26:14:2080:G:O4'	2.20	0.42
1:13:46:G:H2'	1:13:366:C:C5	2.55	0.42
5:42:84:PHE:N	5:42:87:SER:O	2.51	0.42
26:1H:2620:C:OP1	29:21:152:LYS:O	2.37	0.42
34:58:99:LEU:O	34:58:103:VAL:HG23	2.20	0.42
33:61:93:THR:HG22	33:61:119:PRO:HB3	2.01	0.42
1:13:1304:G:N1	1:13:1332:A:OP2	2.48	0.42
26:14:1788:C:C2	26:14:1789:A:C8	3.07	0.42
26:14:511:U:H5''	26:14:512:G:OP2	2.19	0.42
1:13:665:A:C5	1:13:733:A:C5	3.07	0.42
46:D5:48:PHE:CE1	46:D5:52:SER:HA	2.55	0.42
1:1G:1103:C:C2	1:1G:1104:G:C8	3.08	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
28:19:76:PRO:HA	28:19:118:VAL:HG23	2.01	0.42
27:16:17:C:H2'	27:16:18:G:O4'	2.20	0.42
26:14:2459:A:C5	26:14:2460:U:C5	3.08	0.42
26:1H:345:A:N3	26:1H:347:A:N6	2.68	0.42
50:H5:23:LEU:HD11	50:H5:53:LEU:HD13	2.01	0.42
1:1G:32:A:H2'	1:1G:33:A:C8	2.55	0.42
2:1E:109:SER:O	2:1E:112:VAL:HB	2.20	0.42
26:14:2883:A:H5'	26:14:2884:U:H5'	2.02	0.42
41:85:46:ALA:HB1	41:85:50:ARG:HH11	1.85	0.42
7:6E:43:PHE:O	7:6E:46:ALA:HB3	2.20	0.42
26:1H:2456:C:H6	26:1H:2456:C:O5'	2.03	0.42
51:I5:25:TYR:N	51:I5:25:TYR:CD1	2.87	0.42
19:AI:32:LYS:HG2	19:AI:32:LYS:H	1.56	0.42
26:14:442:G:C6	26:14:444:C:N4	2.87	0.42
32:51:94:TYR:HA	32:51:106:THR:O	2.20	0.42
26:1H:2270:G:OP2	58:1H:4056:HOH:O	2.22	0.42
26:14:676:A:H2	26:14:802:A:H61	1.58	0.42
26:1H:654(D):G:H22	26:1H:654(Q):C:H42	1.66	0.42
1:1G:376:G:H1	1:1G:387:U:H3	1.68	0.42
1:1G:192:U:C4'	20:BA:103:GLY:HA2	2.50	0.42
45:C5:97:ARG:HG2	45:C5:102:CYS:O	2.20	0.42
26:14:2495:G:O3'	37:45:81:VAL:HG12	2.20	0.42
34:15:40:PRO:O	41:85:64:ARG:HG2	2.19	0.42
26:1H:1021:A:C3'	26:1H:1021:A:C8	3.01	0.42
26:14:1791:A:H3'	26:14:1792:G:H8	1.84	0.42
1:1G:1349:A:P	9:82:118:LYS:NZ	2.92	0.42
1:13:259:G:H1	1:13:267:C:H42	1.68	0.42
4:3E:9:CYS:O	4:3E:13:ARG:HG3	2.20	0.42
26:1H:247:G:H4'	26:1H:386:G:C5	2.55	0.42
1:1G:1302:U:H5	13:4A:17:VAL:HG21	1.84	0.42
40:B8:16:ARG:HG3	40:B8:18:ASP:OD2	2.20	0.42
26:1H:1817:G:C6	26:1H:1818:U:C5	3.08	0.42
1:13:447:G:C6	1:13:485:G:H1'	2.55	0.42
2:12:165:VAL:HG23	2:12:166:ASP:H	1.85	0.42
26:14:528:A:H2	26:14:2043:C:C5'	2.33	0.42
20:BI:43:LEU:HB2	20:BI:52:ALA:HB2	2.00	0.42
26:1H:634:C:H2'	26:1H:635:C:C6	2.54	0.42
13:4I:27:LYS:HD3	13:4I:31:LYS:NZ	2.34	0.42
1:1G:1329:A:H2'	1:1G:1330:U:O4'	2.20	0.42
1:13:1014:A:N3	1:13:1219:U:H1'	2.35	0.42
24:3K:9:A:H62	24:3K:23:A:H62	1.66	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:13:21:G:H2'	1:13:22:G:C8	2.55	0.42
26:14:28:A:C2	26:14:513:A:C8	3.07	0.42
26:14:839:U:H2'	26:14:840:C:H6	1.84	0.42
26:14:2106:G:C2	26:14:2184:G:C2	3.07	0.42
26:1H:1124:C:H2'	26:1H:1125:G:O4'	2.20	0.42
13:4A:92:HIS:HE2	13:4A:98:VAL:HG21	1.84	0.42
26:1H:197:A:H62	26:1H:2430:A:H2'	1.85	0.42
15:6A:75:PRO:HB2	15:6A:79:ARG:NH2	2.35	0.42
12:3A:32:PHE:HE1	12:3A:86:ARG:HG3	1.85	0.42
10:1I:96:ILE:O	10:1I:97:GLU:HG2	2.19	0.42
26:1H:493:G:H2'	26:1H:494:G:O4'	2.19	0.42
4:32:91:SER:OG	4:32:191:ARG:HG3	2.20	0.42
26:14:1889:A:H2'	26:14:1890:A:O4'	2.20	0.42
26:1H:1094:U:O2	26:1H:1096:A:H5'	2.20	0.42
35:68:22:ILE:HG23	35:68:22:ILE:HD12	1.80	0.42
26:14:565:C:H4'	26:14:1253:A:C6	2.54	0.42
26:1H:470:A:H2'	26:1H:471:A:O4'	2.20	0.42
26:1H:714:U:O2'	26:1H:716:A:N7	2.47	0.42
5:42:28:PHE:CE1	5:42:51:VAL:HG22	2.55	0.42
4:32:107:ARG:HH22	4:32:196:LEU:HD21	1.85	0.42
26:1H:363(F):A:H4'	26:1H:364:C:H5'	2.02	0.42
26:14:1750:G:O2'	26:14:1751:C:H5'	2.20	0.42
28:19:244:ARG:HB2	28:19:245:PRO:HD2	2.02	0.42
26:14:433:C:C4	26:14:434:U:O4	2.73	0.42
1:13:303:A:C5	1:13:304:U:C5	3.07	0.42
42:95:24:LYS:HA	42:95:92:THR:OG1	2.19	0.42
26:14:678:C:H2'	26:14:679:C:C6	2.55	0.42
16:7I:40:ASP:O	16:7I:42:ARG:N	2.53	0.42
11:2I:62:GLN:HB2	11:2I:93:GLN:OE1	2.20	0.42
35:68:9:GLU:O	35:68:83:ALA:HA	2.20	0.42
26:1H:526:A:N3	26:1H:2044:C:H1'	2.35	0.42
49:K8:25:VAL:O	49:K8:29:LYS:HG3	2.20	0.42
27:16:75:G:H21	46:H8:85:HIS:CE1	2.38	0.42
42:D8:2:PHE:O	42:D8:42:GLY:N	2.53	0.42
36:78:81:GLN:OE1	36:78:106:LEU:HA	2.20	0.42
26:1H:2354:G:N2	26:1H:2355:C:C2	2.88	0.42
26:14:244:A:C2	26:14:255:A:C4	3.07	0.42
55:Q8:59:LYS:HG3	55:Q8:59:LYS:HZ2	1.61	0.42
18:9A:41:LYS:O	18:9A:41:LYS:HD3	2.20	0.42
18:9A:51:LEU:HA	18:9A:51:LEU:HD23	1.87	0.42
31:41:62:LEU:HD12	31:41:62:LEU:HA	1.72	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
50:H5:4:LEU:O	50:H5:36:VAL:HA	2.19	0.42
7:6E:65:ALA:HB2	7:6E:128:ALA:HB2	2.00	0.42
26:1H:1615:C:H2'	26:1H:1617:C:C5	2.55	0.42
26:1H:589:C:H2'	26:1H:590:A:C8	2.56	0.41
2:12:213:LEU:O	2:12:213:LEU:HD23	2.20	0.41
1:13:453:A:C4'	16:7I:72:ARG:HB2	2.48	0.41
26:1H:58:G:C2	26:1H:70:G:C6	3.08	0.41
40:B8:88:ILE:HD13	40:B8:91:ARG:NH2	2.34	0.41
29:29:81:ILE:HG23	29:29:81:ILE:HD12	1.81	0.41
27:1J:73:A:C4	27:1J:104:A:C2	3.08	0.41
41:C8:92:ARG:HD3	41:C8:92:ARG:O	2.20	0.41
1:13:1331:G:H5''	1:13:1331:G:H8	1.85	0.41
26:14:2250:G:C8	26:14:2496:C:H5'	2.55	0.41
26:14:2687:U:C4	26:14:2688:U:C5	3.08	0.41
26:14:2002:G:C6	58:14:3900:HOH:O	2.70	0.41
6:5E:97:PHE:O	18:9I:31:LEU:HD23	2.19	0.41
1:13:1126:U:O4	1:13:1127:G:N1	2.53	0.41
49:K8:47:ASN:ND2	49:K8:47:ASN:H	2.17	0.41
49:K8:48:HIS:C	49:K8:48:HIS:ND1	2.73	0.41
26:14:882:G:H2'	26:14:883:G:H8	1.85	0.41
26:1H:242:G:C8	55:Q8:3:LYS:HE3	2.55	0.41
1:1G:553:A:H2'	1:1G:554:C:C6	2.54	0.41
1:13:129:U:N3	1:13:131:C:N4	2.67	0.41
42:95:38:LEU:HD23	42:95:52:VAL:HG21	2.02	0.41
7:6E:95:ARG:HH21	7:6E:99:LEU:HD11	1.85	0.41
3:22:73:PRO:HG3	3:22:105:GLU:HG3	2.01	0.41
26:1H:2816:C:H2'	26:1H:2817:G:C8	2.55	0.41
1:13:198:G:N7	1:13:220:G:N2	2.68	0.41
29:29:62:PRO:C	29:29:64:LYS:H	2.23	0.41
26:14:2115:G:O2'	26:14:2171:A:N6	2.51	0.41
26:14:600:G:N2	26:14:605:C:O3'	2.53	0.41
22:1K:37:MIA:H2'	22:1K:38:A:O4'	2.20	0.41
26:14:2790:A:H4'	26:14:2791:C:OP2	2.20	0.41
1:1G:1478:C:O2'	1:1G:1479:C:H5'	2.20	0.41
10:1A:78:ASN:OD1	10:1A:81:THR:HG23	2.21	0.41
45:C5:64:GLU:HG3	45:C5:64:GLU:H	1.57	0.41
2:1E:226:ARG:HG3	2:1E:227:GLY:N	2.35	0.41
1:1G:109:A:H5'	1:1G:110:C:C5	2.55	0.41
33:69:2:LYS:H	33:69:2:LYS:HG2	1.56	0.41
26:14:2787:C:H4'	29:29:63:LEU:HD21	2.02	0.41
26:1H:1093:G:O2'	26:1H:1099:G:N2	2.52	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:1E:131:PRO:HD2	2:1E:134:GLU:HG3	2.02	0.41
42:D8:91:TYR:HD1	42:D8:91:TYR:H	1.67	0.41
41:C8:28:ARG:O	41:C8:35:ALA:HA	2.20	0.41
7:6E:142:GLU:O	7:6E:146:GLU:HG2	2.20	0.41
1:13:864:A:H3'	1:13:865:A:C8	2.55	0.41
26:14:2490:G:N3	26:14:2490:G:H2'	2.35	0.41
31:49:36:LYS:HG2	31:49:38:VAL:HG23	2.01	0.41
46:D5:178:GLU:HG3	46:D5:179:ASP:H	1.85	0.41
26:14:1742:C:H5'	26:14:1743:G:OP2	2.20	0.41
15:6A:27:VAL:O	15:6A:31:LEU:HB2	2.20	0.41
8:72:31:PHE:HZ	8:72:134:ILE:HD11	1.85	0.41
27:1J:19:G:H2'	27:1J:20:C:O4'	2.20	0.41
45:C5:65:ALA:HA	45:C5:66:PRO:HD3	1.94	0.41
2:1E:77:ALA:O	2:1E:81:VAL:HG12	2.20	0.41
7:62:13:GLN:HA	7:62:14:PRO:HD3	1.74	0.41
6:52:26:ILE:O	6:52:30:LEU:HG	2.20	0.41
26:14:2704:C:H2'	26:14:2705:A:O4'	2.20	0.41
11:2I:66:LEU:HD21	11:2I:97:ALA:O	2.20	0.41
1:1G:693:G:H2'	1:1G:694:A:C8	2.55	0.41
43:A5:51:LEU:HD23	43:A5:51:LEU:HA	1.79	0.41
38:98:54:LEU:HA	38:98:54:LEU:HD12	1.85	0.41
16:7A:25:ARG:HH11	16:7A:25:ARG:HG3	1.85	0.41
47:I8:55:ARG:HG3	47:I8:55:ARG:HH11	1.85	0.41
26:1H:978:G:C2	26:1H:986:C:C2	3.07	0.41
26:1H:2137:C:H2'	26:1H:2138:C:C6	2.55	0.41
9:82:5:TYR:CE2	9:82:7:THR:HG23	2.54	0.41
26:14:993:G:C4	26:14:994:C:H5	2.38	0.41
44:B5:44:GLU:HG2	44:B5:49:VAL:O	2.20	0.41
26:1H:782:A:H5'	26:1H:783:A:C2	2.55	0.41
10:1I:34:VAL:O	10:1I:34:VAL:HG23	2.21	0.41
26:14:2001:A:H2'	26:14:2002:G:C8	2.55	0.41
8:7E:118:VAL:O	8:7E:119:LEU:HD23	2.20	0.41
26:14:2439:A:O2'	26:14:2440:C:OP2	2.31	0.41
28:19:242:ARG:HG3	28:19:246:PRO:HG3	2.02	0.41
49:K8:50:ILE:HD12	49:K8:50:ILE:N	2.35	0.41
1:13:1124:G:N7	1:13:1145:C:H2'	2.35	0.41
1:1G:1347:G:H22	1:1G:1374:A:P	2.43	0.41
4:32:13:ARG:HD2	4:32:38:TYR:O	2.19	0.41
38:98:52:ILE:O	38:98:55:ALA:N	2.53	0.41
50:H5:40:THR:CG2	50:H5:43:ILE:HG12	2.46	0.41
29:29:6:GLY:HA2	29:29:51:PHE:CZ	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:747:U:C6	52:J5:2:ALA:HB3	2.54	0.41
1:1G:1054:C:N4	24:1L:34:G:C4	2.89	0.41
26:14:446:G:H8	58:14:3676:HOH:O	2.03	0.41
13:4A:29:ARG:HB3	13:4A:64:TRP:CZ2	2.55	0.41
18:9A:21:LYS:NZ	18:9A:57:GLY:HA3	2.33	0.41
3:22:22:TRP:HB3	3:22:59:ARG:HB2	2.01	0.41
1:13:142:G:C2	1:13:143:A:N7	2.88	0.41
26:1H:1561:G:O2'	26:1H:1562:A:H5'	2.19	0.41
26:1H:528:A:N1	26:1H:2043:C:H4'	2.34	0.41
1:13:20:U:H2'	1:13:21:G:O4'	2.20	0.41
4:3E:141:ARG:HB2	4:3E:141:ARG:CZ	2.50	0.41
45:G8:9:LYS:HA	45:G8:27:VAL:CG2	2.50	0.41
34:58:35:ARG:HB3	34:58:37:LYS:HG3	2.02	0.41
35:68:47:ILE:HG12	35:68:48:PRO:HD2	2.02	0.41
1:1G:284:G:H2'	1:1G:285:G:C8	2.55	0.41
5:42:145:LYS:HG2	5:42:149:GLU:HG3	2.02	0.41
46:H8:48:PHE:CE1	46:H8:71:VAL:HG11	2.55	0.41
26:1H:479:A:O2'	26:1H:481:G:H2'	2.20	0.41
1:1G:605:U:H2'	1:1G:606:G:O4'	2.20	0.41
1:13:35:G:O2'	12:3I:118:SER:O	2.22	0.41
45:C5:85:VAL:HG23	45:C5:96:ILE:HG22	2.02	0.41
26:14:1184:G:C6	26:14:1185:C:C4	3.09	0.41
17:8A:99:SER:O	17:8A:100:LYS:NZ	2.47	0.41
1:13:380:G:N2	1:13:384:G:C5	2.89	0.41
37:45:58:PHE:HD2	37:45:61:GLY:HA3	1.84	0.41
1:13:924:C:H2'	1:13:925:G:C8	2.54	0.41
26:1H:255:A:H1'	26:1H:384:U:C6	2.55	0.41
4:3E:17:VAL:HG11	4:3E:197:PRO:HB3	2.02	0.41
28:11:72:LYS:HD2	28:11:75:ILE:HD12	2.01	0.41
41:85:14:HIS:O	41:85:18:LEU:HD12	2.19	0.41
1:1G:186(E):C:C2	1:1G:191(C):G:N2	2.88	0.41
8:72:75:ARG:HA	8:72:76:PRO:HD2	1.84	0.41
26:1H:2574:G:O2'	29:21:143:ASN:HB3	2.20	0.41
38:98:86:ARG:HB3	38:98:118:GLU:OE2	2.19	0.41
5:42:39:GLY:O	5:42:69:VAL:N	2.40	0.41
3:2E:73:PRO:O	3:2E:76:VAL:HG13	2.21	0.41
1:1G:1490:C:C4	1:1G:1491:G:N7	2.88	0.41
26:14:1399:C:H2'	26:14:1400:G:H8	1.84	0.41
1:13:1510:U:H2'	1:13:1511:G:C8	2.56	0.41
26:14:17:G:H2'	26:14:18:C:H6	1.84	0.41
26:1H:2241:A:O2'	26:1H:2242:G:H5'	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
40:B8:105:LEU:HD12	40:B8:105:LEU:HA	1.75	0.41
41:C8:9:VAL:HG23	41:C8:9:VAL:H	1.60	0.41
26:14:1801:G:OP2	28:19:154:LYS:NZ	2.47	0.41
29:29:105:THR:HG21	29:29:164:ARG:CZ	2.50	0.41
26:1H:2436:G:C6	26:1H:2437:U:C4	3.09	0.41
26:1H:2468:G:H5'	37:88:120:ILE:HD12	2.02	0.41
1:1G:963:G:H21	10:1A:55:LYS:CE	2.13	0.41
26:14:1614:A:H2	58:14:3515:HOH:O	2.02	0.41
26:1H:1187:G:P	58:1H:3822:HOH:O	2.78	0.41
30:39:29:ASN:HB3	30:39:112:MET:HE1	2.03	0.41
1:13:963:G:C2	10:1I:55:LYS:NZ	2.84	0.41
36:78:2:LYS:HE3	36:78:4:SER:CB	2.43	0.41
1:13:1036:G:H5'	1:13:1037:C:OP2	2.20	0.41
1:1G:1248:A:C6	1:1G:1249:C:N4	2.88	0.41
47:I8:49:LYS:HE2	47:I8:80:HIS:CB	2.50	0.41
47:I8:48:GLY:HA3	47:I8:80:HIS:ND1	2.35	0.41
31:41:67:LYS:CE	51:M8:6:HIS:CE1	3.02	0.41
1:1G:984:C:H2'	1:1G:985:C:C6	2.55	0.41
1:13:659:U:C2	1:13:660:G:C8	3.08	0.41
1:13:1132:C:H2'	1:13:1133:G:O4'	2.20	0.41
26:1H:1323:U:H2'	26:1H:1324:G:H5'	2.01	0.41
26:14:1534:G:H5'	26:14:1535:U:OP2	2.21	0.41
47:I8:14:ARG:HE	47:I8:14:ARG:H	1.68	0.41
1:13:233:C:O2'	1:13:234:C:H5'	2.20	0.41
26:1H:270(T):G:C6	26:1H:270(U):C:C4	3.09	0.41
1:1G:1224:G:N1	1:1G:1322:C:H1'	2.35	0.41
55:Q8:9:GLY:N	55:Q8:12:LYS:HG3	2.31	0.41
4:32:104:VAL:O	4:32:108:LEU:HB2	2.19	0.41
3:22:70:VAL:HG21	3:22:76:VAL:HG11	2.02	0.41
26:1H:1274:A:N1	26:1H:1644:C:O2'	2.40	0.41
7:6E:48:LYS:HD2	7:6E:49:ILE:HD13	2.03	0.41
42:D8:75:PHE:HD1	42:D8:82:ARG:HG3	1.86	0.41
5:4E:136:MET:O	5:4E:140:ARG:NH1	2.53	0.41
26:14:2418:A:OP1	55:M5:29:LYS:NZ	2.52	0.41
30:31:122:LYS:HD2	30:31:191:ARG:HG2	2.02	0.41
17:8A:21:VAL:O	17:8A:41:LYS:HA	2.20	0.41
26:14:2622:C:H5'	29:29:159:HIS:ND1	2.35	0.41
26:14:111:A:C2	26:14:112:U:C2	3.08	0.41
3:22:123:GLN:O	3:22:126:ARG:HB2	2.20	0.41
1:1G:446:G:H2'	1:1G:447:G:O4'	2.20	0.41
1:13:947:G:H2'	1:13:948:C:C6	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:13:501:C:H2'	1:13:502:G:C8	2.56	0.41
26:14:2176:A:H2'	26:14:2177:C:C6	2.55	0.41
4:32:14:ARG:NH1	4:32:14:ARG:HG3	2.35	0.41
19:AA:71:LEU:HA	19:AA:71:LEU:HD23	1.77	0.41
28:19:72:LYS:HB3	28:19:75:ILE:HD12	2.02	0.41
26:14:610:C:H2'	26:14:611:C:H6	1.85	0.41
26:14:370:G:H4'	26:14:371:A:OP2	2.21	0.41
2:12:125:PRO:HA	2:12:127:ILE:HG12	2.02	0.41
26:14:752:A:OP1	54:L5:3:ARG:NH2	2.44	0.41
16:7I:56:ALA:O	16:7I:60:LEU:HD12	2.20	0.41
1:1G:438:G:H5'	4:32:123:HIS:ND1	2.35	0.41
42:D8:35:LEU:HD12	42:D8:37:VAL:O	2.21	0.41
26:1H:2352:A:H2'	26:1H:2353:G:O4'	2.20	0.41
49:K8:41:ILE:HD13	49:K8:41:ILE:O	2.20	0.41
1:13:478:A:O5'	1:13:478:A:H8	2.03	0.41
1:13:397:A:N3	1:13:397:A:H3'	2.36	0.41
31:49:105:LYS:HB2	31:49:105:LYS:HE3	1.92	0.41
26:1H:974(A):C:H4'	26:1H:975:G:C5'	2.50	0.41
1:13:540:G:H2'	1:13:541:G:O4'	2.20	0.41
19:AI:25:LYS:HD3	19:AI:27:GLU:HB2	2.02	0.41
26:14:495:G:H21	43:A5:61:ASN:HD21	1.68	0.41
26:1H:972:G:OP2	26:1H:973:A:O2'	2.28	0.41
26:1H:1359:A:N3	26:1H:1359:A:O4'	2.53	0.41
26:14:1022:G:N2	26:14:1142(A):A:H2	2.16	0.41
26:14:2552:U:H2'	26:14:2554:U:OP2	2.20	0.41
55:Q8:46:ARG:HB2	55:Q8:46:ARG:NH1	2.35	0.41
26:1H:1526:G:H2'	26:1H:1527:G:O4'	2.21	0.41
11:2A:54:ARG:O	11:2A:57:THR:OG1	2.39	0.41
26:14:2131:G:C5'	26:14:2158:A:H61	2.33	0.41
52:N8:40:LYS:HE2	52:N8:47:PRO:HD2	2.02	0.41
1:13:812:C:H2'	1:13:812:C:H6	1.52	0.41
40:75:80:SER:HA	40:75:81:PRO:HD3	1.95	0.41
26:14:2376:A:H2'	26:14:2377:A:C8	2.55	0.41
1:1G:430:A:C4	1:1G:431:A:C8	3.08	0.41
1:1G:468:A:O2'	16:7A:81:ARG:HA	2.19	0.41
12:3I:7:ILE:HD12	12:3I:7:ILE:HG23	1.82	0.41
1:1G:500:G:H2'	1:1G:501:C:H6	1.84	0.41
40:75:23:ARG:HG3	40:75:120:ARG:HH12	1.84	0.41
1:1G:406:G:C2	1:1G:407:G:N7	2.88	0.41
26:14:1777:U:O2'	26:14:1778:U:H5'	2.19	0.41
26:1H:2300:G:H1	26:1H:2316:C:H42	1.67	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:303:U:H2'	26:14:304:G:O4'	2.21	0.41
24:3L:33:U:OP2	24:3L:33:U:H2'	2.21	0.41
55:Q8:34:TRP:CG	55:Q8:35:GLN:N	2.86	0.41
1:13:789:U:H5	1:13:792:A:OP2	2.03	0.41
1:1G:1121:U:H2'	1:1G:1122:U:O4'	2.20	0.41
1:1G:994:A:H2'	1:1G:995:C:C6	2.55	0.41
26:1H:754:C:H2'	26:1H:755:C:H6	1.82	0.41
26:1H:2404:C:N4	26:1H:2405:G:C6	2.88	0.41
26:14:959:A:C6	26:14:960:A:N1	2.87	0.41
1:1G:1084:G:C8	1:1G:1085:U:C6	3.08	0.41
12:3I:102:ARG:HE	12:3I:102:ARG:HB3	1.58	0.41
26:14:1165:U:H2'	26:14:1166:C:C6	2.55	0.41
35:68:64:ARG:HG2	35:68:79:PHE:CD2	2.56	0.41
1:1G:188:U:O2'	1:1G:189:U:H5'	2.21	0.41
26:14:776:G:H4'	26:14:777:A:O5'	2.19	0.41
1:1G:255:G:C2	1:1G:272:C:C2	3.07	0.41
12:3A:10:LEU:HD13	17:8A:32:TYR:CZ	2.55	0.41
26:14:1743:G:C2	26:14:1746:G:C8	3.08	0.41
33:61:64:GLU:OE1	33:61:67:ARG:NH1	2.52	0.41
42:D8:98:GLU:HG2	42:D8:99:ILE:N	2.34	0.41
9:82:26:VAL:HG13	9:82:61:ALA:O	2.20	0.41
45:C5:6:HIS:CD2	45:C5:7:VAL:HG13	2.56	0.41
28:11:232:PRO:HA	58:11:408:HOH:O	2.20	0.41
20:BI:51:GLU:O	20:BI:54:LYS:HB3	2.20	0.41
26:1H:2031:A:C6	26:1H:2498:C:H1'	2.56	0.41
15:6I:8:LYS:O	15:6I:12:ILE:HG13	2.20	0.41
43:E8:39:THR:HG22	43:E8:44:ALA:HB2	2.02	0.41
42:D8:53:GLU:HG2	42:D8:54:GLY:N	2.36	0.41
46:H8:111:VAL:O	46:H8:114:GLY:HA2	2.20	0.41
23:2K:59:A:H4'	23:2K:60:A:OP1	2.21	0.41
1:13:1513:A:H2'	1:13:1514:C:C6	2.55	0.41
26:14:1799:G:O6	28:19:178:PRO:HD2	2.21	0.41
42:95:72:VAL:HG13	42:95:72:VAL:O	2.21	0.41
12:3A:60:LEU:HD13	12:3A:60:LEU:HA	1.74	0.41
32:59:152:ARG:HD2	32:59:153:LYS:HG3	2.02	0.41
1:13:16:A:N1	1:13:919:A:H2	2.18	0.41
26:14:723:G:H2'	26:14:724:U:O4'	2.20	0.41
43:E8:64:MET:O	43:E8:65:LEU:HB2	2.20	0.41
36:78:19:VAL:CB	36:78:20:GLY:HA2	2.51	0.41
2:12:5:ILE:HB	2:12:221:LEU:HD21	2.03	0.41
15:6I:26:GLU:H	15:6I:26:GLU:HG2	1.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
16:7I:72:ARG:O	16:7I:75:ARG:HB3	2.20	0.41
22:1K:12:U:O2	22:1K:24:G:N2	2.54	0.41
33:69:128:LEU:O	33:69:137:PRO:HA	2.21	0.41
1:1G:521:G:O5'	12:3A:73:GLU:HG2	2.20	0.41
26:1H:1533:C:H2'	26:1H:1534:G:C4	2.55	0.41
26:1H:1534:G:N2	26:1H:1538:G:N2	2.68	0.41
30:39:122:LYS:HB2	30:39:123:LEU:H	1.52	0.41
26:1H:1022:G:N2	26:1H:1142(A):A:N1	2.67	0.41
37:88:59:ARG:C	37:88:61:GLY:N	2.73	0.41
1:1G:474:G:C2	1:1G:475:G:C5	3.09	0.41
1:13:1133:G:H2'	1:13:1134:G:H8	1.83	0.41
1:1G:279:A:H4'	1:1G:280:C:H5''	2.03	0.41
4:32:173:TRP:HA	4:32:187:ARG:HG2	2.02	0.41
26:14:1011:G:C2	26:14:1013:C:C2	3.09	0.41
16:7A:52:ASP:OD2	16:7A:55:ARG:HG3	2.21	0.41
26:14:2378:A:O2'	39:65:21:THR:HG21	2.20	0.41
26:1H:2316:C:H1'	31:41:128:ARG:NH2	2.35	0.41
1:1G:374:A:H2'	1:1G:374:A:N3	2.35	0.41
3:22:11:ARG:O	3:22:14:ILE:O	2.39	0.41
28:19:11:PRO:O	28:19:12:SER:OG	2.25	0.41
17:8A:82:MET:O	17:8A:86:GLU:N	2.43	0.41
26:1H:1643:G:C6	26:1H:1644:C:C4	3.08	0.41
5:42:136:MET:HG2	5:42:136:MET:H	1.58	0.41
39:65:26:LEU:O	39:65:88:ASP:HB2	2.20	0.41
1:13:48:C:H6	1:13:365:U:O4	2.03	0.41
20:BI:38:LYS:O	20:BI:41:ILE:HG13	2.21	0.41
26:14:1053:C:N4	26:14:1106:G:H1	2.17	0.41
1:1G:918:A:H2'	1:1G:919:A:O4'	2.21	0.41
26:14:2772:C:H2'	26:14:2773:C:C6	2.55	0.41
32:59:73:ALA:O	32:59:76:VAL:HB	2.21	0.41
26:14:2756:U:C2	26:14:2757:A:N7	2.88	0.41
13:4I:56:LEU:HA	13:4I:56:LEU:HD22	1.84	0.41
27:1J:78:A:C2	27:1J:99:A:C4	3.09	0.41
45:C5:43:ASN:HB2	45:C5:62:GLU:O	2.20	0.41
34:15:135:PRO:HB2	34:15:137:LYS:NZ	2.36	0.41
1:1G:867:G:O2'	1:1G:868:C:H5'	2.21	0.41
38:55:104:ARG:HB2	38:55:107:ASP:OD1	2.21	0.41
19:AA:11:VAL:HB	19:AA:12:ASP:H	1.78	0.41
8:7E:75:ARG:HA	8:7E:76:PRO:HD2	1.81	0.41
13:4A:50:GLU:O	13:4A:53:VAL:HB	2.20	0.41
26:14:923:C:H2'	26:14:924:C:C6	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:2E:21:ARG:NH2	3:2E:56:ASP:OD1	2.53	0.41
54:P8:30:VAL:O	54:P8:34:ARG:HG3	2.21	0.41
1:1G:1205:U:H1'	3:22:195:VAL:CG2	2.51	0.41
1:1G:1137:C:H5''	1:1G:1138:G:OP1	2.21	0.41
26:14:1310:G:OP2	54:L5:9:ARG:NE	2.53	0.41
38:98:65:LEU:HD12	38:98:65:LEU:HA	1.68	0.41
51:I5:8:LYS:HD3	51:I5:8:LYS:HA	1.55	0.41
20:BA:58:LYS:O	20:BA:58:LYS:HD3	2.21	0.41
26:14:1033:U:H3'	26:14:1033:U:H6	1.84	0.41
35:68:106:LEU:HD23	35:68:106:LEU:HA	1.77	0.41
40:75:82:LEU:HD12	40:75:82:LEU:H	1.84	0.41
6:5E:10:LEU:HD13	6:5E:61:LEU:HD13	2.02	0.41
26:1H:311:A:C6	26:1H:328:U:C4	3.09	0.41
43:A5:59:VAL:HG12	43:A5:60:ASN:N	2.34	0.41
26:1H:1250:G:OP2	36:78:18:ARG:NH1	2.54	0.41
27:1J:109:G:C6	27:1J:110:G:C5	3.09	0.41
24:3L:36:A:C8	25:4L:14:A:N6	2.88	0.41
1:1G:533:A:P	58:1G:1728:HOH:O	2.78	0.41
39:A8:34:HIS:O	39:A8:97:ARG:NH2	2.54	0.41
1:13:515:G:N2	1:13:537:G:C4	2.89	0.41
2:12:153:ARG:HG3	2:12:154:LEU:HD23	2.01	0.41
46:H8:152:ALA:HB3	46:H8:167:PRO:HA	2.03	0.41
26:1H:444:C:H4'	30:31:49:ALA:HB2	2.03	0.41
26:14:309:G:O3'	45:C5:18:GLY:HA2	2.20	0.41
26:14:2128:C:C4	26:14:2129:C:C4	3.08	0.41
51:M8:59:PHE:O	51:M8:63:TYR:HB2	2.20	0.41
26:14:273(C):C:N3	26:14:363(C):G:N2	2.59	0.41
24:3K:7:A:H2	24:3K:67:C:O2	2.04	0.41
34:58:22:THR:HB	34:58:25:ARG:HB2	2.03	0.41
9:82:118:LYS:O	9:82:119:ALA:HB3	2.20	0.41
1:13:76:G:H1'	1:13:95:G:N2	2.35	0.41
29:29:33:VAL:CG2	29:29:47:VAL:HG13	2.47	0.41
26:14:882:G:H1	26:14:894:C:N4	2.18	0.41
1:1G:1015:A:C6	1:1G:1016:A:C5	3.09	0.41
1:1G:1014:A:H5'	19:AA:15:LEU:HG	2.03	0.41
51:I5:37:SER:C	51:I5:39:CYS:H	2.24	0.41
51:I5:37:SER:OG	51:I5:38:LYS:N	2.53	0.41
1:13:1131:G:H2'	1:13:1132:C:H6	1.85	0.41
38:98:63:ARG:HB2	38:98:80:PHE:CE2	2.53	0.41
27:16:7:G:H5''	27:16:7:G:C8	2.54	0.41
1:13:411:A:N9	1:13:413:G:H1'	2.35	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:12:21:ARG:HB2	2:12:22:LYS:H	1.78	0.41
3:22:50:ALA:HB1	3:22:70:VAL:HG11	2.02	0.41
1:1G:627:G:C2	1:1G:628:G:C8	3.09	0.41
1:13:143:A:H5''	1:13:144:G:H5'	2.02	0.41
31:41:101:ILE:O	31:41:105:LYS:HE2	2.21	0.41
26:1H:2098:U:H2'	26:1H:2099:U:O4'	2.20	0.41
26:14:1385:G:C4	26:14:1386:C:C5	3.09	0.41
22:1K:37:MIA:H121	22:1K:38:A:C2	2.55	0.41
26:14:1174:A:N6	26:14:1176:G:O2'	2.54	0.41
46:D5:108:PRO:HB2	46:D5:110:GLY:O	2.20	0.41
7:62:69:VAL:HG22	7:62:135:VAL:HG22	2.01	0.41
37:88:33:GLY:HA2	37:88:105:GLU:HA	2.03	0.41
37:88:11:LYS:NZ	37:88:86:GLY:HA2	2.36	0.41
33:69:8:PRO:HD3	33:69:15:VAL:HG22	2.01	0.41
26:1H:722:A:H2'	26:1H:723:G:C8	2.56	0.41
26:14:2698:U:H2'	26:14:2699:C:C6	2.56	0.41
20:BA:50:GLU:N	20:BA:100:ILE:HG12	2.36	0.41
1:13:971:G:N1	1:13:1363:A:OP2	2.50	0.41
4:32:148:VAL:HG12	4:32:152:SER:OG	2.21	0.41
1:1G:440:A:H8	1:1G:440:A:OP2	2.04	0.41
27:16:88:C:H2'	27:16:89:G:O4'	2.20	0.41
1:1G:1291:G:H2'	1:1G:1292:U:C6	2.55	0.41
46:D5:127:LYS:HB3	46:D5:127:LYS:HE2	1.68	0.41
26:1H:2365:G:H4'	47:I8:60:PHE:CZ	2.55	0.41
27:16:116:G:H2'	27:16:117:G:O4'	2.21	0.41
26:1H:2319:G:H4'	26:1H:2320:A:OP1	2.21	0.41
31:41:31:VAL:HA	31:41:32:PRO:HD3	1.92	0.41
26:14:1137:G:O2'	26:14:2039:C:H5'	2.20	0.41
6:5E:5:GLU:HA	6:5E:63:TYR:O	2.20	0.41
1:1G:664:G:P	18:9A:64:ARG:HH21	2.44	0.41
24:1L:4:C:O2	24:1L:70:G:N2	2.53	0.41
26:1H:2003:G:H2'	26:1H:2004:G:O5'	2.21	0.41
2:12:187:LEU:HA	2:12:201:ILE:O	2.21	0.41
45:C5:44:ILE:HG13	45:C5:45:VAL:N	2.36	0.41
1:13:1396:A:O4'	1:13:1398:A:H1'	2.20	0.41
3:2E:188:LEU:HD13	3:2E:195:VAL:HG11	2.02	0.41
32:51:23:ARG:NH1	32:51:25:LYS:HG3	2.35	0.41
26:14:1016:G:C6	26:14:1017:G:C5	3.08	0.41
1:1G:382:A:H2'	1:1G:383:A:H8	1.86	0.41
26:1H:744:G:P	29:21:132:HIS:HD1	2.43	0.41
26:14:404:C:O2'	26:14:405:U:OP2	2.35	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:58:121:LYS:HB3	34:58:123:TYR:CE1	2.56	0.41
26:14:1759:A:H4'	26:14:2715:C:O4'	2.20	0.41
33:69:52:ARG:HA	33:69:55:ALA:HB3	2.02	0.41
38:98:4:LEU:HA	38:98:4:LEU:HD13	1.38	0.41
37:88:45:GLN:H	37:88:45:GLN:CD	2.23	0.41
27:1J:60:C:H2'	27:1J:61:G:H8	1.85	0.41
2:1E:28:PHE:O	2:1E:32:ILE:HG22	2.20	0.41
26:1H:1783:A:P	58:1H:3626:HOH:O	2.79	0.41
55:Q8:28:GLY:HA3	55:Q8:29:LYS:HA	1.09	0.41
7:6E:28:ASN:O	7:6E:31:MET:HB3	2.21	0.41
1:13:1368:G:C5'	9:8E:112:LYS:HB3	2.38	0.41
26:1H:973:A:O4'	26:1H:1188:U:C6	2.74	0.41
1:1G:533:A:C5	1:1G:536:C:C4	3.09	0.41
26:14:1141:U:H3'	34:15:63:THR:HG21	2.02	0.41
26:1H:1479:G:H2'	26:1H:1480:G:H8	1.83	0.41
55:Q8:58:ILE:HG21	55:Q8:58:ILE:HD13	1.76	0.41
26:1H:2061:G:H2'	26:1H:2501:C:O2'	2.20	0.41
27:1J:27:C:O3'	39:65:36:TYR:OH	2.32	0.41
4:32:22:LYS:N	4:32:26:CYS:SG	2.91	0.41
33:69:128:LEU:HD13	33:69:128:LEU:HA	1.73	0.41
26:1H:996:A:O2'	41:C8:92:ARG:HG3	2.21	0.41
34:58:96:GLU:HB2	34:58:122:VAL:HG12	2.02	0.41
1:1G:1286:A:C8	1:1G:1286:A:C3'	3.00	0.41
1:1G:861:G:H21	1:1G:872:A:H2	1.68	0.41
1:1G:862:C:C2'	1:1G:863:U:H5'	2.51	0.41
1:1G:345:C:OP2	40:75:39:ARG:NH2	2.54	0.41
33:61:135:GLU:HB2	33:61:136:VAL:H	1.73	0.41
14:5I:53:LEU:HB3	14:5I:56:VAL:HG21	2.03	0.41
31:41:101:ILE:HG13	51:M8:25:TYR:O	2.21	0.41
26:1H:993:G:H5''	41:C8:50:ARG:NH2	2.36	0.41
1:1G:338:A:H2	1:1G:351:G:H22	1.68	0.41
31:49:130:ASN:HB3	31:49:160:VAL:HA	2.02	0.41
28:19:239:ARG:CZ	58:19:302:HOH:O	2.68	0.41
26:14:265:A:C8	26:14:266:G:H1'	2.56	0.41
13:4A:92:HIS:CD2	13:4A:98:VAL:HG11	2.54	0.41
17:8A:59:ILE:HG22	17:8A:71:PHE:CD2	2.56	0.41
26:1H:869:G:H2'	26:1H:870:A:O4'	2.21	0.41
26:1H:1824:G:OP1	28:11:52:ARG:HD3	2.21	0.41
26:1H:997:G:OP1	41:C8:93:LYS:N	2.51	0.41
26:1H:2792:G:C6	26:1H:2805:G:N1	2.88	0.41
38:98:38:VAL:HG22	38:98:112:ALA:HB2	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1G:341:C:H2'	1:1G:342:C:H6	1.86	0.41
41:C8:88:ILE:HG22	41:C8:90:VAL:HB	2.01	0.41
1:13:498:A:H4'	1:13:500:G:OP1	2.19	0.41
43:E8:95:ILE:HD13	43:E8:95:ILE:HG21	1.65	0.41
26:1H:1470:G:H5''	26:1H:1471:A:OP1	2.21	0.41
26:14:557:U:H2'	26:14:558:G:H8	1.86	0.41
18:9I:53:ARG:NE	18:9I:59:SER:O	2.50	0.41
23:2K:19:G:C2	23:2K:59:A:C5	3.09	0.41
32:59:23:ARG:HA	32:59:36:PRO:HA	2.02	0.41
1:13:66:G:O4'	1:13:173:U:C4	2.74	0.41
46:D5:23:LYS:HB3	46:D5:38:TYR:CD2	2.55	0.41
26:14:1104:C:H2'	26:14:1105:U:C6	2.55	0.41
26:14:301:G:C4	26:14:302:C:C5	3.09	0.41
32:59:7:LEU:N	32:59:8:PRO:HD2	2.34	0.41
34:15:118:LYS:O	34:15:121:LYS:NZ	2.47	0.41
32:51:56:SER:OG	32:51:58:GLU:HG2	2.19	0.41
7:6E:27:ILE:HD12	7:6E:40:ALA:HA	2.02	0.41
17:8I:9:VAL:HG21	17:8I:84:LEU:HB3	2.02	0.41
26:1H:2356:C:C5	26:1H:2357:U:C4	3.09	0.41
26:14:1860:G:H8	26:14:1860:G:O5'	2.04	0.41
31:49:128:ARG:HD3	31:49:128:ARG:HA	1.79	0.41
3:2E:154:SER:HA	3:2E:165:THR:HA	2.03	0.41
5:4E:96:PRO:HA	5:4E:117:ASP:OD2	2.21	0.41
26:1H:1523:U:H2'	26:1H:1524:G:O4'	2.21	0.41
36:35:97:PRO:HG3	36:35:112:LEU:CB	2.50	0.41
26:1H:574:C:H4'	26:1H:575:A:O5'	2.21	0.41
27:1J:109:G:C5	27:1J:110:G:N7	2.88	0.41
26:14:1020:A:H61	26:14:1141:U:HO2'	1.68	0.41
26:1H:878:A:C2	26:1H:879:G:C5	3.08	0.41
36:78:50:ARG:CG	36:78:50:ARG:HH21	2.27	0.41
27:1J:66:A:C2	27:1J:108:C:C4	3.09	0.41
36:78:144:GLU:O	36:78:144:GLU:OE1	2.39	0.41
26:1H:654(P):G:H2'	26:1H:654(Q):C:C6	2.56	0.41
26:1H:654(B):C:H42	26:1H:654(S):G:H1	1.68	0.41
1:13:57:G:C5	1:13:58:C:C4	3.08	0.41
1:13:874:G:C6	1:13:875:C:C4	3.09	0.41
2:12:131:PRO:HG2	2:12:134:GLU:HG3	2.02	0.41
1:1G:683:G:H2'	1:1G:684:A:C8	2.55	0.41
26:14:603:A:H8	26:14:604:G:H1'	1.86	0.41
1:1G:660:G:OP2	15:6A:5:LYS:HD3	2.21	0.41
26:14:1328:G:H2'	26:14:1330:C:C4	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
31:41:124:SER:HB2	31:41:131:TYR:CE1	2.56	0.41
26:14:2393:A:C2'	26:14:2394:C:H5'	2.51	0.41
26:14:1639:U:OP2	58:14:3529:HOH:O	2.22	0.41
13:4A:94:ARG:NH2	19:AA:78:ARG:HH12	2.18	0.41
26:14:1993:U:H4'	29:29:128:SER:CB	2.50	0.41
1:1G:827:U:H5''	1:1G:828:A:OP2	2.21	0.41
26:14:1149:G:C2	26:14:1150:C:N3	2.88	0.41
37:88:72:LYS:HA	37:88:73:PRO:HD3	1.84	0.41
9:8E:89:ASN:C	9:8E:91:ASP:H	2.24	0.41
26:14:867:C:C5	26:14:868:U:C5	3.09	0.41
26:14:288:C:H2'	26:14:289:A:C8	2.55	0.41
26:1H:729:G:C6	28:11:208:LYS:HB2	2.55	0.41
1:1G:559:A:H4'	1:1G:560:U:C5'	2.51	0.41
1:13:1314:C:N3	1:13:1315:U:C4	2.89	0.41
1:1G:865:A:H5'	1:1G:1078:U:C5	2.56	0.41
44:F8:11:PRO:HG2	44:F8:13:LEU:HD21	2.03	0.41
26:14:1599:C:H2'	26:14:1600:C:C6	2.56	0.41
1:1G:991:U:OP2	1:1G:991:U:H6	2.04	0.41
1:13:46:G:H2'	1:13:366:C:H5	1.85	0.41
26:1H:1831:G:C4	26:1H:1975:G:N2	2.89	0.41
45:C5:85:VAL:CG2	45:C5:98:VAL:HB	2.50	0.41
1:13:946:A:H2'	1:13:947:G:C8	2.56	0.41
1:13:683:G:H2'	1:13:684:A:H8	1.84	0.41
26:14:1695:G:H2'	26:14:1696:G:O4'	2.20	0.41
40:75:57:PHE:HA	40:75:79:HIS:HD2	1.86	0.41
54:P8:26:GLY:O	54:P8:30:VAL:HG23	2.21	0.41
36:35:124:LYS:HA	36:35:143:GLY:O	2.21	0.41
26:1H:1906:G:O2'	26:1H:1907:G:H5'	2.21	0.41
26:1H:545:G:H2'	26:1H:546:C:H5''	2.02	0.41
45:C5:99:CYS:SG	45:C5:100:ALA:N	2.93	0.41
26:1H:731:C:C2	26:1H:732:C:C5	3.09	0.41
26:14:2329:G:N2	47:E5:41:ARG:HB3	2.36	0.41
1:1G:105:G:C6	1:1G:106:C:C4	3.08	0.41
26:14:2422:A:O2'	26:14:2423:U:H5''	2.20	0.41
33:61:35:LEU:O	33:61:36:ALA:HB2	2.21	0.41
16:7I:8:ARG:O	16:7I:9:PHE:HD1	2.03	0.41
26:1H:1853:A:H2'	26:1H:1854:A:C8	2.56	0.41
50:H5:18:ASP:O	50:H5:21:ALA:N	2.51	0.41
1:1G:1260:C:H6	1:1G:1260:C:H3'	1.85	0.41
2:1E:55:PHE:HD1	2:1E:55:PHE:HA	1.68	0.41
43:E8:90:ARG:HB3	43:E8:90:ARG:HE	1.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
43:E8:96:ILE:HD12	43:E8:96:ILE:HG23	1.75	0.41
47:I8:32:ARG:HB3	47:I8:32:ARG:HE	1.52	0.41
16:7I:55:ARG:HA	16:7I:55:ARG:HD2	1.71	0.41
10:1A:35:SER:HB3	10:1A:73:ASP:HB2	2.01	0.41
26:14:2569:G:C2	26:14:2570:G:C8	3.09	0.41
26:1H:618:G:H2'	26:1H:618(A):C:H6	1.85	0.41
26:1H:811:U:H2'	36:78:21:ARG:HA	2.03	0.41
26:1H:607:U:O2	26:1H:621:A:N1	2.54	0.41
26:14:1614:A:H61	43:A5:88:ARG:H	1.67	0.41
1:13:1399:C:C2	1:13:1401:G:C5	3.09	0.41
44:B5:31:HIS:HA	44:B5:32:PRO:HD3	1.89	0.41
7:6E:113:GLU:HG3	7:6E:119:ARG:HA	2.03	0.41
26:14:1021:A:C8	26:14:1021:A:C3'	3.03	0.41
1:13:452:A:H62	1:13:480:U:H3	1.68	0.41
26:1H:2592:G:C6	26:1H:2593:U:C4	3.08	0.41
26:14:1226:G:P	42:95:85:LYS:HD3	2.60	0.41
55:Q8:53:PRO:HB3	55:Q8:56:GLU:H	1.86	0.41
24:3K:21:A:H2'	24:3K:22:G:O4'	2.21	0.41
22:1K:14:A:C5	22:1K:22:G:N2	2.89	0.41
1:1G:1436:U:H2'	1:1G:1437:C:O4'	2.21	0.41
27:1J:40:U:O2'	27:1J:45:A:N6	2.44	0.41
26:14:2134:A:C8	26:14:2158:A:N7	2.89	0.41
1:1G:1207:G:H2'	1:1G:1208:C:H6	1.85	0.41
1:13:187:C:H1'	1:13:191(A):G:N2	2.36	0.41
26:14:2410:G:H2'	26:14:2411:A:O4'	2.21	0.41
34:15:39:ARG:HH11	34:15:48:MET:HE2	1.86	0.41
28:19:31:LYS:O	28:19:32:SER:OG	2.37	0.41
28:19:70:TRP:CD1	28:19:71:ASP:N	2.89	0.41
26:1H:266:G:N2	26:1H:427:U:H1'	2.35	0.41
32:59:140:LYS:HB3	32:59:140:LYS:HE2	1.81	0.41
26:1H:1176:G:H5'	26:1H:1177:A:C8	2.55	0.41
31:41:37:VAL:HG22	31:41:159:VAL:HG12	2.03	0.41
26:1H:2157:G:HO2'	26:1H:2158:A:P	2.43	0.41
4:3E:13:ARG:HD3	4:3E:36:ARG:O	2.21	0.41
26:14:94:G:H21	49:G5:47:ASN:HD22	1.67	0.41
1:1G:502:G:H2'	1:1G:503:C:O4'	2.20	0.41
26:14:2017:U:P	58:14:3835:HOH:O	2.78	0.41
26:1H:887:A:H5'	26:1H:888:C:OP1	2.21	0.41
26:1H:2286:A:H8	53:O8:37:ARG:HH11	1.69	0.41
40:B8:19:LEU:HA	40:B8:20:PRO:HD3	1.90	0.41
26:1H:662:G:C5'	36:78:15:ARG:H	2.30	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:1340:U:H4'	26:14:1394:U:O2'	2.21	0.41
29:21:50:GLY:CA	29:21:77:ILE:HG12	2.51	0.41
30:39:129:PHE:HA	30:39:142:TRP:NE1	2.35	0.41
36:78:64:LYS:HE3	55:Q8:12:LYS:HD2	2.03	0.41
2:1E:69:LEU:HD13	2:1E:91:PRO:HB2	2.02	0.41
33:69:76:THR:HG21	33:69:140:LEU:HA	2.02	0.41
5:4E:37:ARG:HH12	5:4E:111:GLU:HG2	1.86	0.41
1:1G:115:G:H4'	1:1G:116:A:O5'	2.20	0.41
28:11:108:PRO:HG3	28:11:143:HIS:HE1	1.81	0.41
26:1H:2244:U:H6	26:1H:2244:U:O5'	2.03	0.41
28:11:68:LYS:HB3	28:11:70:TRP:CZ2	2.55	0.41
26:1H:1420:U:HO2'	26:1H:1421:G:P	2.43	0.41
26:14:729:G:OP2	28:19:13:ARG:NH1	2.52	0.41
33:61:130:TYR:O	33:61:135:GLU:HB2	2.21	0.41
7:62:143:ARG:NH1	24:3L:42:C:O5'	2.54	0.41
44:B5:18:TYR:C	44:B5:20:GLY:H	2.23	0.41
8:7E:9:MET:O	8:7E:12:ARG:N	2.54	0.41
26:14:2488:A:H8	26:14:2488:A:O5'	2.04	0.41
46:H8:80:ARG:HG2	46:H8:80:ARG:H	1.48	0.41
29:29:96:PHE:HD2	29:29:182:LEU:HD21	1.85	0.41
28:11:36:PRO:HA	28:11:61:LEU:CD1	2.51	0.41
26:1H:1429:G:C4	26:1H:1568:G:C2	3.08	0.41
32:51:98:LEU:HD13	32:51:125:VAL:CG2	2.50	0.41
26:14:1317:A:H2'	26:14:1318:C:C6	2.56	0.41
14:5I:3:ARG:HH21	14:5I:6:LEU:HG	1.86	0.41
32:59:9:ILE:HD12	32:59:49:VAL:HB	2.02	0.41
26:1H:1029:A:N1	26:1H:2465:C:O2'	2.48	0.41
1:13:458:C:N3	1:13:475:G:N2	2.69	0.41
26:14:193:U:H5	58:14:3625:HOH:O	2.03	0.41
11:2I:57:THR:HG22	11:2I:59:TYR:N	2.36	0.41
29:21:4:ILE:HG12	29:21:5:LEU:H	1.85	0.41
9:82:46:ALA:O	9:82:49:PRO:HD2	2.21	0.41
30:39:21:ALA:O	30:39:23:ASP:N	2.54	0.41
10:1I:5:ARG:NH1	10:1I:99:LYS:HD2	2.36	0.41
10:1I:89:ASP:C	10:1I:91:PRO:HD3	2.41	0.41
37:45:69:PHE:HA	37:45:70:PRO:HD2	1.81	0.41
26:14:491:G:H2'	26:14:492:A:C8	2.56	0.41
26:14:146:G:H2'	26:14:147:U:O4'	2.21	0.41
1:1G:776:G:N2	1:1G:802:A:OP2	2.43	0.41
26:14:2575:C:H2'	26:14:2578:G:O6	2.21	0.41
38:55:52:ILE:O	38:55:55:ALA:N	2.53	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
11:2I:73:MET:CG	11:2I:103:LEU:HD23	2.50	0.41
26:1H:182:A:H2'	26:1H:183:C:C6	2.56	0.41
44:F8:41:ASN:OD1	44:F8:41:ASN:N	2.54	0.41
44:F8:21:PHE:CD1	44:F8:26:TYR:CD1	3.08	0.41
26:1H:216:A:H2'	26:1H:217:G:C8	2.56	0.41
26:14:1100:C:H2'	26:14:1101:U:C6	2.56	0.41
48:F5:73:LEU:HB3	48:F5:90:ILE:HG23	2.03	0.41
12:3A:110:VAL:HG23	12:3A:120:TYR:HB3	2.03	0.41
17:8I:18:THR:HG23	17:8I:69:LYS:HD2	2.03	0.41
26:1H:1475:G:H2'	26:1H:1476:C:C6	2.56	0.41
39:65:24:LEU:HB2	39:65:85:VAL:CG1	2.50	0.41
38:98:118:GLU:HA	38:98:118:GLU:OE1	2.20	0.41
6:5E:61:LEU:HB3	6:5E:63:TYR:HE2	1.86	0.41
26:14:2780:G:OP1	34:15:118:LYS:HE2	2.21	0.41
28:19:232:PRO:HA	58:19:301:HOH:O	2.21	0.41
15:6I:36:ILE:HA	15:6I:59:MET:CE	2.51	0.41
26:14:49:A:H4'	26:14:50:U:H5''	2.02	0.41
9:82:15:ALA:HB2	9:82:65:VAL:HB	2.03	0.41
47:E5:36:ILE:HD11	47:E5:39:ARG:HG2	2.03	0.41
26:1H:2590:A:OP2	28:11:238:GLY:HA3	2.20	0.41
26:1H:1007:C:OP2	26:1H:1008:C:O2'	2.35	0.41
41:85:43:GLY:HA3	42:95:73:SER:OG	2.21	0.41
1:1G:1267:C:O2	21:1B:20:LYS:HD2	2.20	0.41
26:14:774:A:H2	26:14:787:U:HO2'	1.60	0.41
42:D8:34:GLU:HG3	42:D8:56:SER:OG	2.19	0.41
1:13:1152:A:C6	1:13:1153:C:C4	3.09	0.41
26:14:608:A:H2'	26:14:609:A:C8	2.56	0.41
4:32:139:ARG:HG3	4:32:139:ARG:HH11	1.86	0.41
26:1H:1776:G:H2'	26:1H:1776:G:N3	2.36	0.41
26:14:1568:G:P	28:19:63:ARG:HH12	2.42	0.41
1:13:542:G:H5'	4:3E:41:GLY:HA3	2.03	0.41
26:1H:2450:A:C2	26:1H:2451:A:C4	3.09	0.41
26:14:2822:G:O2'	26:14:2824:C:OP2	2.31	0.41
37:88:69:PHE:HA	37:88:70:PRO:HD2	1.90	0.41
26:14:2484:G:C2	26:14:2485:G:C8	3.09	0.41
50:L8:2:PRO:HB2	50:L8:3:ARG:H	1.66	0.41
26:14:1910:G:H1	26:14:1920:C:H42	1.68	0.41
19:AI:11:VAL:HG13	19:AI:16:LEU:HD22	2.03	0.41
26:1H:2724:C:OP1	29:21:118:LYS:HE3	2.21	0.41
26:1H:2663:G:C6	26:1H:2664:G:C4	3.09	0.41
17:8I:46:ASP:OD2	17:8I:49:GLU:HG3	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:1H:2443:C:O2'	26:1H:2444:G:H5'	2.20	0.41
26:1H:988:A:P	50:L8:11:SER:HB2	2.61	0.41
29:21:181:LEU:HD12	29:21:181:LEU:HA	1.77	0.41
50:L8:38:GLU:H	50:L8:38:GLU:CD	2.24	0.41
32:51:153:LYS:CE	32:51:153:LYS:H	2.34	0.41
13:4A:47:ASP:OD1	13:4A:47:ASP:N	2.54	0.41
5:4E:151:LEU:HD23	5:4E:151:LEU:HA	1.93	0.41
46:H8:70:LEU:HD23	46:H8:70:LEU:HA	1.97	0.41
1:13:246:A:C2	1:13:282:A:C5	3.09	0.41
26:1H:2396:G:H5''	48:J8:25:LYS:NZ	2.36	0.41
54:L5:19:ARG:HG2	54:L5:19:ARG:NH1	2.34	0.41
26:1H:1728:G:C2	26:1H:1730:U:OP2	2.74	0.41
44:F8:1:MET:C	44:F8:3:THR:N	2.73	0.41
26:1H:2035:G:H4'	26:1H:2036:C:OP2	2.21	0.41
1:13:1182:G:C8	1:13:1182:G:O5'	2.74	0.41
1:13:1181:G:O2'	1:13:1184:G:H5'	2.21	0.41
36:35:50:ARG:CB	36:35:50:ARG:HH11	2.34	0.41
26:1H:449:A:C6	26:1H:450:G:C5	3.09	0.41
1:13:76:G:H2'	1:13:77:C:H5'	2.02	0.41
26:1H:1177:A:OP1	26:1H:1178:C:H5	2.03	0.41
1:1G:1013:G:O2'	1:1G:1014:A:N7	2.38	0.41
34:58:94:HIS:C	34:58:95:PRO:O	2.59	0.41
26:1H:880:G:H2'	26:1H:881:G:H8	1.83	0.41
36:78:24:GLY:C	36:78:26:GLY:N	2.74	0.41
35:68:78:ARG:HH21	40:B8:103:ARG:NH2	2.18	0.41
38:98:87:TYR:CD1	38:98:90:ARG:HD2	2.53	0.41
4:32:59:ARG:NH2	4:32:66:ARG:NH1	2.70	0.41
12:3I:8:ASN:HA	12:3I:11:VAL:HG23	2.03	0.41
3:22:47:LEU:HD22	3:22:47:LEU:HA	1.84	0.41
39:A8:27:SER:HA	39:A8:88:ASP:CB	2.48	0.41
1:1G:1326:C:H2'	1:1G:1327:C:H6	1.83	0.41
15:6A:61:GLY:O	15:6A:65:ARG:HG3	2.21	0.41
1:1G:1227:A:C8	1:1G:1227:A:H3'	2.56	0.41
1:1G:823:G:H2'	1:1G:824:C:C6	2.56	0.41
27:16:15:A:H1'	27:16:109:G:N9	2.35	0.41
5:42:40:ARG:HA	5:42:67:VAL:O	2.20	0.41
39:65:102:ALA:O	39:65:105:ALA:N	2.54	0.41
26:1H:1181:C:O2'	26:1H:1182:A:H5'	2.21	0.41
14:5I:6:LEU:HB3	14:5I:23:ARG:HH22	1.86	0.41
10:1I:6:ILE:HD11	10:1I:72:VAL:HB	2.03	0.41
1:13:580:U:H3	1:13:761:G:H1	1.69	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
48:F5:56:GLN:NE2	48:F5:83:GLU:HA	2.35	0.41
1:13:805:C:O2'	1:13:806:C:H5'	2.21	0.41
39:A8:63:THR:O	39:A8:66:ALA:HB3	2.20	0.41
23:2L:54:G:O2'	23:2L:55:5MU:H5''	2.21	0.41
26:1H:2864:G:H2'	26:1H:2865:U:H6	1.85	0.41
1:1G:735:C:H2'	1:1G:736:C:H6	1.85	0.41
1:13:622:A:H2'	1:13:623:C:O4'	2.21	0.41
11:2A:59:TYR:CZ	11:2A:63:LEU:HD12	2.56	0.41
1:1G:440:A:H3'	1:1G:442:C:C5	2.56	0.41
1:13:1303:C:N4	1:13:1304:G:C6	2.89	0.41
8:7E:108:GLY:HA3	8:7E:138:TRP:HB3	2.03	0.41
38:98:29:LEU:HA	38:98:29:LEU:HD12	1.68	0.41
4:32:94:LEU:HA	4:32:97:LEU:HD12	2.01	0.41
1:13:595:G:H22	1:13:643:C:H41	1.69	0.41
26:14:2370:G:C6	26:14:2371:G:C6	3.08	0.41
7:6E:27:ILE:CD1	7:6E:40:ALA:HA	2.51	0.41
26:14:909:A:C8	26:14:912:C:N4	2.89	0.41
1:13:936:C:H2'	1:13:937:A:H5'	2.03	0.41
1:13:1057:G:C5	1:13:1204:A:C2	3.09	0.41
32:59:92:ILE:HB	32:59:93:GLY:H	1.64	0.41
41:C8:27:LEU:O	41:C8:30:LYS:N	2.54	0.41
54:P8:24:THR:HA	54:P8:25:PRO:HD2	1.96	0.41
26:1H:2626:C:H2'	26:1H:2627:G:O4'	2.21	0.41
47:E5:54:GLY:O	47:E5:57:PHE:N	2.52	0.41
1:1G:772:U:H2'	1:1G:773:G:O4'	2.20	0.41
26:14:1514:U:H2'	26:14:1515:C:C6	2.56	0.41
26:14:1444(A):A:N3	26:14:1444(A):A:H2'	2.36	0.41
17:8A:74:LEU:HA	17:8A:74:LEU:HD22	1.85	0.41
14:5I:44:LEU:HD12	14:5I:44:LEU:O	2.21	0.41
45:G8:44:ILE:HG13	45:G8:44:ILE:H	1.55	0.41
7:6E:10:ARG:HE	7:6E:10:ARG:HB2	1.61	0.41
40:75:95:ARG:HA	40:75:95:ARG:HD2	1.92	0.41
26:14:1447:G:H1'	26:14:1545(A):A:H1'	2.02	0.41
30:39:155:LEU:HB2	30:39:189:THR:HG21	2.03	0.41
31:41:11:TYR:O	31:41:16:ARG:HG3	2.21	0.41
48:J8:25:LYS:HB3	48:J8:25:LYS:HE3	1.90	0.40
26:14:1899:G:O2'	26:14:1900:A:H5''	2.22	0.40
1:1G:589:C:C2	1:1G:650:G:N2	2.84	0.40
26:1H:71:A:OP1	26:1H:72:U:H2'	2.21	0.40
48:J8:41:ARG:HH11	48:J8:41:ARG:CG	2.22	0.40
24:3K:7:A:H2	24:3K:67:C:C2	2.39	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:58:1:MET:HE1	41:C8:95:LEU:HD11	2.03	0.40
1:13:1004:A:H1'	1:13:1036:G:N2	2.34	0.40
47:I8:49:LYS:HB2	47:I8:49:LYS:HE3	1.89	0.40
31:41:67:LYS:NZ	51:M8:6:HIS:CE1	2.89	0.40
45:G8:95:LYS:HE2	45:G8:97:ARG:HH22	1.86	0.40
12:3I:53:ARG:HB2	12:3I:93:LEU:HD11	2.03	0.40
26:14:2017:U:O2	52:J5:10:LYS:HB2	2.21	0.40
26:1H:1454:U:H5'	38:98:63:ARG:CZ	2.51	0.40
10:1A:12:ASP:HB3	10:1A:15:THR:HG23	2.03	0.40
26:1H:2261:C:O4'	26:1H:2388:A:H1'	2.21	0.40
30:31:37:VAL:HG21	36:78:6:LEU:HD21	2.03	0.40
32:59:41:MET:HB3	32:59:42:ARG:H	1.71	0.40
1:1G:625:G:H2'	1:1G:626:U:H6	1.85	0.40
4:3E:89:THR:H	4:3E:92:VAL:CG2	2.34	0.40
29:29:64:LYS:HB3	29:29:65:GLY:H	1.59	0.40
26:14:2875:C:OP1	40:75:3:ARG:NH1	2.54	0.40
1:1G:23:C:H5	1:1G:561:U:O4	2.03	0.40
53:K5:36:LEU:HB3	53:K5:50:ARG:HH11	1.86	0.40
7:62:18:TYR:HB3	7:62:59:LEU:CD1	2.51	0.40
13:4I:39:ILE:HD12	13:4I:56:LEU:HD23	2.03	0.40
26:1H:934:G:H2'	26:1H:935:C:C6	2.56	0.40
29:29:112:GLY:O	29:29:159:HIS:HA	2.22	0.40
28:19:176:ARG:HH11	28:19:176:ARG:HG2	1.85	0.40
1:1G:1194:U:H4'	5:42:22:GLY:O	2.21	0.40
26:14:2735:G:H2'	26:14:2736:G:H8	1.86	0.40
23:2K:24:C:H2'	23:2K:25:U:C6	2.56	0.40
1:13:339:C:OP2	35:68:97:ARG:NH1	2.44	0.40
36:35:65:ARG:HB3	36:35:65:ARG:HH11	1.86	0.40
2:1E:165:VAL:HG23	2:1E:166:ASP:N	2.36	0.40
1:13:114:U:O2'	1:13:115:G:H5'	2.20	0.40
26:14:1751:C:H2'	26:14:1752:C:C6	2.56	0.40
26:1H:2356:C:H2'	26:1H:2357:U:O4'	2.21	0.40
31:41:12:TYR:HA	31:41:16:ARG:HG3	2.04	0.40
24:3K:51:U:N3	24:3K:63:G:O6	2.54	0.40
39:65:59:LYS:HZ3	39:65:61:ASN:HA	1.87	0.40
26:14:1436:G:O2'	26:14:1477:A:H4'	2.21	0.40
28:19:136:ILE:HG22	28:19:140:THR:OG1	2.21	0.40
26:1H:1545(A):A:N7	26:1H:1546:C:O2	2.54	0.40
17:8A:3:LYS:HB3	17:8A:61:GLU:HB3	2.03	0.40
26:1H:458:G:O2'	54:P8:39:ARG:HD3	2.21	0.40
26:14:1638:C:H4'	26:14:2710:C:O2	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:13:1521:G:H2'	1:13:1522:U:C6	2.55	0.40
26:1H:1680:U:H2'	26:1H:1681:G:O4'	2.21	0.40
26:14:2187:G:C6	26:14:2188:C:N3	2.90	0.40
3:22:28:GLN:CD	3:22:28:GLN:H	2.24	0.40
7:62:89:MET:HB3	7:62:89:MET:HE2	1.87	0.40
31:49:88:ILE:HD12	31:49:88:ILE:HA	1.88	0.40
46:D5:123:ASP:N	46:D5:123:ASP:OD1	2.53	0.40
26:14:357:A:C2	26:14:358:U:C2	3.09	0.40
38:98:103:ARG:CZ	38:98:110:PRO:HD3	2.50	0.40
41:C8:78:THR:O	41:C8:81:HIS:N	2.54	0.40
26:1H:575:A:OP2	26:1H:2055:C:N4	2.54	0.40
1:1G:1149:C:P	9:82:9:ARG:HH11	2.44	0.40
1:13:926:G:C6	1:13:1505:G:C5	3.09	0.40
26:14:2356:C:H4'	47:E5:20:ARG:HG3	2.02	0.40
46:H8:126:VAL:HA	46:H8:164:ALA:H	1.85	0.40
44:B5:9:LEU:HA	49:G5:36:ARG:HH21	1.86	0.40
26:14:2304:G:H5'	31:49:132:ASN:HB2	2.03	0.40
2:1E:100:GLY:HA2	2:1E:103:THR:OG1	2.21	0.40
48:J8:73:LEU:HA	48:J8:73:LEU:HD23	1.87	0.40
1:1G:191(F):U:H2'	1:1G:191:G:H8	1.86	0.40
26:1H:1539:G:C2	26:1H:1540:G:C5	3.10	0.40
26:14:259:G:N2	26:14:621:A:H8	2.15	0.40
1:13:607:A:C2	16:7I:31:LYS:HG3	2.55	0.40
3:2E:40:ARG:HG3	3:2E:40:ARG:NH1	2.34	0.40
26:1H:1331:A:O2'	26:1H:1332:G:C8	2.70	0.40
24:3L:5:G:H2'	24:3L:6:G:C8	2.56	0.40
26:14:2294:C:P	39:65:89:ARG:HH22	2.44	0.40
29:29:5:LEU:HD12	29:29:197:ILE:HG22	2.03	0.40
23:2K:20:G:C4	23:2K:58:A:C2	3.09	0.40
26:14:2299:G:C2	26:14:2318:G:H8	2.39	0.40
1:13:1285:A:H4'	1:13:1286:A:C5'	2.52	0.40
39:65:29:PHE:O	39:65:35:ILE:HD12	2.21	0.40
23:2L:57:C:H2'	23:2L:58:A:C8	2.56	0.40
35:68:75:SER:OG	40:B8:74:ARG:NH2	2.54	0.40
1:13:1179:A:O3'	9:8E:103:THR:HG23	2.21	0.40
44:F8:57:LEU:HD23	44:F8:57:LEU:N	2.35	0.40
26:1H:2689:U:C6	26:1H:2689:U:H5'	2.56	0.40
1:13:1342:C:H2'	1:13:1343:G:C8	2.56	0.40
20:BI:14:LYS:HG2	20:BI:14:LYS:O	2.20	0.40
33:69:3:VAL:HG12	33:69:38:LEU:HA	2.03	0.40
26:14:236:C:H2'	26:14:237:C:C6	2.56	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:1907:G:C2	26:14:1924:C:O2	2.75	0.40
26:14:2:G:H22	26:14:2900:A:H61	1.67	0.40
26:1H:251:A:C5	26:1H:252:G:H1'	2.56	0.40
5:42:93:PRO:HG2	8:72:105:ARG:HE	1.86	0.40
45:G8:40:GLU:HA	45:G8:41:GLY:HA2	1.82	0.40
1:1G:1307:U:O5'	1:1G:1307:U:H6	2.04	0.40
37:45:54:MET:HE1	37:45:104:PHE:HB3	2.03	0.40
1:1G:254:G:OP1	17:8A:67:LYS:O	2.40	0.40
1:13:1442:G:C5	1:13:1446:A:C6	3.09	0.40
9:82:53:VAL:HG11	9:82:92:TYR:CZ	2.56	0.40
4:3E:116:GLN:HE22	4:3E:157:LEU:HD11	1.87	0.40
26:14:1337:G:H2'	26:14:1338:G:C8	2.54	0.40
37:88:25:ASP:H	37:88:102:VAL:HG22	1.87	0.40
4:3E:8:VAL:C	4:3E:10:ARG:N	2.75	0.40
1:1G:763:G:H2'	1:1G:764:C:H6	1.86	0.40
26:14:2619:C:H2'	26:14:2620:C:H6	1.86	0.40
45:G8:43:ASN:CG	45:G8:67:LEU:HD11	2.42	0.40
30:39:161:GLU:O	30:39:165:ARG:HG3	2.22	0.40
41:C8:25:TRP:O	41:C8:28:ARG:HB2	2.21	0.40
26:14:142:G:H2'	26:14:143:C:H6	1.86	0.40
26:1H:2402:C:H5	26:1H:2415:G:H22	1.70	0.40
17:8I:9:VAL:O	17:8I:21:VAL:HA	2.22	0.40
1:13:929:G:C6	1:13:930:C:C4	3.10	0.40
29:29:134:ILE:O	29:29:135:HIS:C	2.59	0.40
1:1G:1313:U:OP1	19:AA:7:LYS:N	2.54	0.40
39:A8:74:ALA:HB1	39:A8:107:GLU:O	2.21	0.40
40:75:106:SER:HA	40:75:110:ILE:HD12	2.03	0.40
26:14:2607:G:H2'	26:14:2608:G:O4'	2.22	0.40
30:39:132:VAL:HG22	30:39:133:ASN:H	1.85	0.40
1:13:1414:U:H2'	1:13:1415:G:H8	1.86	0.40
29:29:108:SER:OG	29:29:163:GLU:HG2	2.22	0.40
29:29:14:ILE:HB	40:75:14:TYR:CE2	2.56	0.40
27:1J:51:G:C6	27:1J:52:A:H2	2.39	0.40
16:7A:82:GLN:HG2	16:7A:82:GLN:H	1.61	0.40
2:12:55:PHE:HA	2:12:55:PHE:HD1	1.78	0.40
26:1H:2745:C:H4'	32:51:142:GLY:O	2.21	0.40
26:14:2101:G:H2'	26:14:2102:U:O4'	2.22	0.40
51:M8:37:SER:O	51:M8:41:PRO:HD2	2.22	0.40
9:82:16:ARG:O	9:82:63:ILE:HG23	2.20	0.40
26:14:71:A:C2	44:B5:31:HIS:NE2	2.78	0.40
26:1H:1727:U:H2'	26:1H:1728:G:O4'	2.22	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:13:537:G:H2'	1:13:538:G:H8	1.87	0.40
42:95:8:GLY:O	42:95:10:LYS:HE3	2.21	0.40
26:1H:1582:C:HO2'	26:1H:1586:A:H8	1.68	0.40
26:14:310:A:OP1	45:C5:17:SER:O	2.40	0.40
26:1H:2126:A:H62	26:1H:2163:C:H1'	1.86	0.40
26:14:2392:A:C2	26:14:2424:C:N4	2.79	0.40
1:1G:1068:G:N3	1:1G:1191:A:C2	2.89	0.40
42:95:76:LYS:NZ	42:95:82:ARG:HH21	2.19	0.40
28:19:93:ALA:HB3	28:19:105:ILE:HG22	2.04	0.40
1:13:632:A:C8	1:13:633:G:C4	3.09	0.40
38:55:37:THR:OG1	38:55:39:PRO:HD2	2.20	0.40
1:13:607:A:O2'	1:13:608:A:H5'	2.20	0.40
51:M8:12:ALA:C	51:M8:24:THR:HG21	2.41	0.40
26:14:2307:G:HO2'	26:14:2308:G:P	2.35	0.40
29:29:47:VAL:O	29:29:47:VAL:HG12	2.21	0.40
26:1H:412:A:H2'	26:1H:412:A:N3	2.37	0.40
1:1G:545:C:H2'	1:1G:546:G:O4'	2.22	0.40
4:32:60:GLU:HG2	4:32:202:LEU:HB2	2.03	0.40
26:1H:459:U:H2'	26:1H:460:A:H8	1.86	0.40
26:14:1639:U:P	58:14:3529:HOH:O	2.78	0.40
38:98:63:ARG:NH2	38:98:80:PHE:HD2	2.19	0.40
26:1H:2231:C:H2'	26:1H:2232:U:O4'	2.21	0.40
28:11:131:LEU:HB2	28:11:136:ILE:CD1	2.51	0.40
1:13:342:C:N3	1:13:347:G:N2	2.62	0.40
33:69:76:THR:HG23	33:69:77:LEU:N	2.35	0.40
12:3A:78:GLN:HG3	12:3A:81:SER:CB	2.51	0.40
2:1E:11:LEU:HB3	2:1E:213:LEU:HD11	2.02	0.40
1:13:123:C:OP1	1:13:312:C:H5'	2.21	0.40
26:1H:65:C:H2'	26:1H:66:C:H6	1.86	0.40
26:14:1795:C:H2'	26:14:1796:U:C6	2.56	0.40
1:13:247:G:C2	1:13:248:C:C6	3.09	0.40
1:1G:485:G:HO2'	1:1G:486:U:P	2.43	0.40
26:1H:1337:G:C2	26:1H:1338:G:C4	3.10	0.40
39:65:66:ALA:HA	39:65:69:VAL:CG1	2.52	0.40
26:1H:806:C:OP2	36:78:41:ARG:HD3	2.21	0.40
26:1H:1635:G:C4	26:1H:1636:C:C5	3.10	0.40
11:2I:85:ARG:HD3	11:2I:113:PRO:HD3	2.01	0.40
26:1H:253:C:H2'	26:1H:254:G:O4'	2.22	0.40
44:B5:27:THR:HG22	44:B5:80:ILE:HG22	2.02	0.40
26:14:221:A:C4	26:14:266:G:N7	2.89	0.40
10:1A:65:LEU:HD12	14:5A:55:GLY:O	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1G:994:A:H2'	1:1G:995:C:H6	1.86	0.40
26:14:2784:C:H2'	26:14:2785:C:C6	2.56	0.40
26:14:1657:C:H2'	26:14:1658:C:H6	1.86	0.40
38:55:79:LEU:HA	38:55:83:ILE:HB	2.02	0.40
36:78:85:LEU:HA	36:78:88:LEU:HD22	2.02	0.40
4:3E:19:LEU:HB3	4:3E:21:LEU:HD21	2.02	0.40
28:19:267:SER:C	28:19:269:PHE:H	2.24	0.40
53:O8:27:LYS:HZ2	53:O8:27:LYS:H	1.68	0.40
51:M8:16:CYS:HG	51:M8:36:CYS:HG	1.68	0.40
26:1H:784:A:C8	26:1H:792:G:C5	3.10	0.40
8:7E:39:LEU:CD1	8:7E:111:ILE:HD11	2.50	0.40
1:13:610:G:H2'	1:13:611:A:O4'	2.21	0.40
26:14:1625:C:H2'	26:14:1626:G:H5'	2.04	0.40
1:13:1350:A:C6	1:13:1351:U:N3	2.90	0.40
1:13:644:G:H2'	1:13:645:C:O4'	2.22	0.40
1:1G:32:A:C2	1:1G:33:A:C4	3.10	0.40
49:K8:29:LYS:HG2	49:K8:57:ILE:HD13	2.04	0.40
8:72:31:PHE:CE2	8:72:35:ILE:HD11	2.56	0.40
34:58:121:LYS:HB3	34:58:123:TYR:HE1	1.86	0.40
54:L5:29:LYS:HA	54:L5:32:LYS:HB2	2.03	0.40
26:14:1032:A:H2	26:14:1122:G:H1	1.69	0.40
1:13:329:A:C5	1:13:332:G:C6	3.09	0.40
1:1G:266:G:H2'	1:1G:266:G:N3	2.37	0.40
26:1H:1161:C:C6	26:1H:1161:C:H3'	2.56	0.40
2:1E:149:LEU:HD23	2:1E:149:LEU:HA	1.91	0.40
31:49:143:GLU:OE2	31:49:143:GLU:N	2.54	0.40
11:2I:107:SER:O	11:2I:108:ILE:HG13	2.21	0.40
1:13:1432:G:N2	1:13:1468:A:OP2	2.55	0.40
1:1G:352:C:P	58:1G:1717:HOH:O	2.79	0.40
40:B8:64:ARG:HH11	40:B8:64:ARG:HD3	1.76	0.40
26:14:1774:C:O5'	26:14:1774:C:H6	2.05	0.40
26:14:1287:A:H5''	26:14:1288:U:OP2	2.21	0.40
26:1H:512:G:N7	58:1H:4190:HOH:O	2.54	0.40
26:1H:2593:U:O2'	26:1H:2594:C:H5'	2.21	0.40
1:13:1024:G:H4'	1:13:1024:G:OP1	2.20	0.40
26:1H:445:C:OP1	41:C8:2:PRO:HA	2.22	0.40
1:13:1180:A:H5''	1:13:1181:G:OP1	2.21	0.40
27:1J:38:C:O4'	39:65:95:HIS:CE1	2.75	0.40
26:1H:2377:A:H2'	26:1H:2378:A:C8	2.56	0.40
24:3K:6:G:H22	24:3K:67:C:N4	2.19	0.40
50:L8:50:VAL:CG2	50:L8:54:VAL:HG11	2.51	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:13:51:A:C6	1:13:353:A:C2	3.10	0.40
1:1G:413:G:H2'	1:1G:428:G:H22	1.84	0.40
51:M8:13:ARG:CA	51:M8:24:THR:HG21	2.51	0.40
1:13:277:C:H2'	1:13:278:G:C8	2.55	0.40
26:14:2388:A:H2'	26:14:2389:G:H5'	2.02	0.40
1:1G:769:G:H4'	1:1G:1513:A:H4'	2.04	0.40
1:1G:1171:G:H2'	1:1G:1172:C:H6	1.84	0.40
1:13:1064:G:H4'	1:13:1065:U:OP1	2.21	0.40
26:14:2844:G:C6	26:14:2845:G:C4	3.09	0.40
15:6I:6:GLU:CD	15:6I:6:GLU:N	2.74	0.40
2:1E:162:ILE:HD13	2:1E:183:PRO:O	2.21	0.40
23:2K:69:C:H2'	23:2K:70:C:C6	2.56	0.40
16:7A:55:ARG:HE	16:7A:55:ARG:HA	1.87	0.40
46:H8:7:ALA:HB3	46:H8:61:LEU:CB	2.50	0.40
1:1G:1330:U:H5'	13:4A:24:GLY:H	1.86	0.40
8:7E:120:THR:H	8:7E:123:GLU:HB2	1.87	0.40
1:1G:1158:C:N3	1:1G:1160:G:C8	2.89	0.40
1:1G:1227:A:H8	1:1G:1227:A:H3'	1.87	0.40
52:J5:36:CYS:HG	52:J5:49:CYS:CB	2.34	0.40
39:65:3:ARG:NH2	39:65:4:LEU:HD12	2.36	0.40
26:1H:536:A:OP1	41:C8:53:ARG:NH1	2.54	0.40
20:BI:14:LYS:HE2	20:BI:14:LYS:HB3	1.78	0.40
11:2I:41:THR:HG21	11:2I:71:LYS:HB3	2.04	0.40
26:1H:2562:U:C1'	35:68:23:ARG:HD3	2.52	0.40
7:62:65:ALA:HB3	7:62:124:LEU:HD22	2.03	0.40
1:1G:265:G:H5'	17:8A:64:PRO:O	2.21	0.40
1:1G:1478:C:H2'	1:1G:1479:C:C6	2.56	0.40
26:1H:686:G:H1	54:P8:16:HIS:CD2	2.40	0.40
18:9I:52:PRO:HB2	18:9I:54:ARG:HG2	2.03	0.40
8:7E:121:ASP:O	8:7E:125:ARG:HB2	2.21	0.40
27:16:60:C:C2	27:16:61:G:C8	3.09	0.40
1:13:560:U:H4'	1:13:561:U:H5''	2.02	0.40
26:1H:775:G:O5'	26:1H:777:A:H1'	2.21	0.40
26:14:458:G:O2'	54:L5:39:ARG:HD3	2.22	0.40
1:13:339:C:H2'	1:13:340:U:H6	1.86	0.40
26:1H:671:C:O2'	26:1H:672:C:H5'	2.21	0.40
48:F5:91:LYS:HE2	48:F5:91:LYS:HB2	1.54	0.40
48:J8:83:GLU:CG	48:J8:85:LEU:HB2	2.51	0.40
41:C8:75:ASN:HB2	41:C8:78:THR:OG1	2.21	0.40
1:1G:730:G:C5	1:1G:731:G:H1'	2.56	0.40
32:59:26:VAL:CG1	32:59:33:LEU:H	2.34	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:2L:73:A:C6	23:2L:74:A:C6	3.10	0.40
1:13:1234:C:O2'	1:13:1235:U:H5'	2.22	0.40
1:13:989:C:H42	1:13:1216:G:H1	1.68	0.40
51:I5:10:VAL:HA	51:I5:11:PRO:HD2	1.82	0.40
24:3L:12:U:C4	24:3L:13:C:C4	3.09	0.40
7:62:60:LYS:HD2	7:62:60:LYS:HA	1.88	0.40
5:4E:101:ILE:HA	5:4E:101:ILE:HD13	1.79	0.40
18:9I:35:ARG:HB2	18:9I:35:ARG:HE	1.64	0.40
34:15:99:LEU:HA	34:15:99:LEU:HD23	1.95	0.40
48:J8:50:ARG:HE	48:J8:50:ARG:HB3	1.76	0.40
26:14:270(D):C:H2'	26:14:270(E):G:C8	2.56	0.40
1:1G:926:G:C6	1:1G:1505:G:C5	3.09	0.40
45:G8:85:VAL:HG23	45:G8:96:ILE:O	2.21	0.40
26:14:1111:A:C4'	32:59:3:ARG:HH11	2.33	0.40
1:13:1022:G:H2'	1:13:1023:G:C8	2.57	0.40
1:13:973:G:H4'	10:1I:54:PHE:O	2.22	0.40
19:AI:40:ILE:O	19:AI:41:VAL:HG22	2.22	0.40
1:1G:674:G:N2	1:1G:717:C:O2	2.55	0.40
15:6A:58:MET:O	15:6A:62:GLN:N	2.37	0.40
1:1G:1288:A:H1'	1:1G:1352:C:O2'	2.21	0.40
5:42:101:ILE:HD13	5:42:101:ILE:H	1.87	0.40
1:13:975:A:C4'	1:13:976:G:H5''	2.46	0.40
26:1H:1668:A:C5	26:1H:1674:G:C5	3.10	0.40
26:1H:1050:A:C8	26:1H:2751:G:C5	3.09	0.40
26:1H:1194:A:OP2	26:1H:1194:A:C8	2.67	0.40
1:1G:1096:C:H2'	1:1G:1097:C:C6	2.57	0.40
26:14:1639:U:C2'	26:14:1640:C:H5'	2.51	0.40
26:1H:2729:G:H2'	26:1H:2730:C:O4'	2.21	0.40
47:E5:24:LYS:O	47:E5:25:ARG:HD3	2.21	0.40
8:7E:41:ARG:NH1	8:7E:123:GLU:OE1	2.54	0.40
26:14:869:G:N2	26:14:870:A:H1'	2.37	0.40
26:1H:2801:A:H5'	26:1H:2895:U:O2'	2.22	0.40
8:72:1:MET:N	8:72:1:MET:SD	2.75	0.40
1:13:628:G:C2	1:13:629:G:C4	3.10	0.40
26:14:1026:U:H5''	26:14:1026:U:H6	1.87	0.40
26:14:945:A:C4	26:14:2448:A:C2	3.08	0.40
42:D8:72:VAL:O	42:D8:72:VAL:HG23	2.22	0.40
2:1E:7:VAL:HG21	2:1E:217:ARG:NH1	2.37	0.40
1:1G:791:G:N1	1:1G:792:A:N6	2.70	0.40
3:2E:124:ILE:HG12	3:2E:130:VAL:HG22	2.04	0.40
1:13:1079:G:H2'	1:13:1080:A:C8	2.57	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:52:33:TYR:CZ	6:52:78:GLU:HG3	2.56	0.40
1:13:1446:A:O2'	40:B8:125:ARG:NH2	2.55	0.40
1:13:407:G:O2'	4:3E:116:GLN:HG3	2.21	0.40
26:14:30:G:C5	26:14:31:C:C4	3.09	0.40
26:1H:1215:G:C5	26:1H:1216:G:C8	3.09	0.40
26:1H:280:C:C2	26:1H:361:G:N2	2.90	0.40
50:H5:13:ILE:HD12	50:H5:13:ILE:H	1.87	0.40
26:14:128:C:H2'	26:14:129:C:C6	2.56	0.40
38:98:8:ARG:NH1	38:98:39:PRO:HB3	2.36	0.40
46:D5:94:GLU:HA	46:D5:95:PRO:HD3	1.86	0.40
46:D5:54:HIS:HB3	46:D5:101:PRO:HG3	2.03	0.40
30:39:64:ILE:HG13	30:39:65:TRP:NE1	2.36	0.40
26:1H:270(X):G:C6	26:1H:270(Y):G:C6	3.09	0.40
33:61:29:TYR:HD2	33:61:30:LEU:HD23	1.87	0.40
26:14:191:A:H1'	26:14:679:C:H1'	2.02	0.40
26:14:2489:G:C6	26:14:2490:G:N7	2.89	0.40
26:14:1469:A:H2'	26:14:1470:G:O4'	2.22	0.40
38:98:96:ARG:NH2	38:98:117:VAL:HG23	2.36	0.40
43:E8:74:ALA:HA	43:E8:104:THR:O	2.21	0.40
46:D5:170:THR:C	46:D5:172:ALA:H	2.24	0.40
26:1H:1918:A:N3	26:1H:1919:A:N6	2.67	0.40
31:41:59:GLU:O	31:41:63:ILE:HG23	2.21	0.40
3:2E:131:ARG:HG2	3:2E:166:GLU:OE2	2.22	0.40
26:1H:1413:G:N2	26:1H:1589:C:O2	2.44	0.40
40:B8:78:LEU:HD12	40:B8:79:HIS:CD2	2.57	0.40
26:1H:2248:C:C5	26:1H:2249:U:C4	3.09	0.40
26:1H:1682:G:C6	26:1H:1683:C:C4	3.10	0.40
32:51:72:ILE:O	32:51:76:VAL:HG23	2.22	0.40
20:BA:59:ALA:HB3	20:BA:84:LEU:HD11	2.03	0.40
26:1H:222:A:H8	26:1H:222:A:H2'	1.80	0.40
44:F8:28:PHE:N	44:F8:28:PHE:CD1	2.89	0.40
1:13:1171:G:O5'	1:13:1171:G:H8	2.04	0.40
26:1H:2263:C:H2'	26:1H:2264:C:H6	1.86	0.40
26:14:2859:G:H3'	26:14:2859:G:C8	2.57	0.40
30:31:8:GLN:CD	30:31:8:GLN:H	2.24	0.40
24:3L:1:G:H8	24:3L:1:G:HO5'	1.70	0.40
8:7E:98:LYS:H	8:7E:98:LYS:HG3	1.74	0.40
4:32:138:TYR:C	4:32:138:TYR:CD2	2.94	0.40
13:4I:16:ASP:N	13:4I:16:ASP:OD1	2.55	0.40
26:1H:2367:G:H2'	26:1H:2368:C:C6	2.56	0.40
26:1H:2555:U:H5''	26:1H:2556:C:OP2	2.22	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:1H:2137:C:OP1	1:1G:999:U:O2'[4_555]	2.14	0.06

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	12	235/256 (92%)	198 (84%)	34 (14%)	3 (1%)	15	48
2	1E	235/256 (92%)	199 (85%)	30 (13%)	6 (3%)	7	29
3	22	204/239 (85%)	178 (87%)	26 (13%)	0	100	100
3	2E	203/239 (85%)	186 (92%)	14 (7%)	3 (2%)	13	44
4	32	206/209 (99%)	179 (87%)	27 (13%)	0	100	100
4	3E	206/209 (99%)	189 (92%)	14 (7%)	3 (2%)	13	44
5	42	149/162 (92%)	139 (93%)	9 (6%)	1 (1%)	26	64
5	4E	149/162 (92%)	139 (93%)	9 (6%)	1 (1%)	26	64
6	52	99/101 (98%)	95 (96%)	4 (4%)	0	100	100
6	5E	99/101 (98%)	94 (95%)	5 (5%)	0	100	100
7	62	153/156 (98%)	143 (94%)	10 (6%)	0	100	100
7	6E	153/156 (98%)	142 (93%)	11 (7%)	0	100	100
8	72	136/138 (99%)	125 (92%)	11 (8%)	0	100	100
8	7E	136/138 (99%)	126 (93%)	9 (7%)	1 (1%)	26	64
9	82	125/128 (98%)	116 (93%)	9 (7%)	0	100	100
9	8E	125/128 (98%)	110 (88%)	15 (12%)	0	100	100
10	1A	97/105 (92%)	88 (91%)	9 (9%)	0	100	100
10	1I	97/105 (92%)	88 (91%)	9 (9%)	0	100	100
11	2A	114/129 (88%)	103 (90%)	10 (9%)	1 (1%)	21	58

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
11	2I	114/129 (88%)	101 (89%)	12 (10%)	1 (1%)	21	58
12	3A	123/132 (93%)	101 (82%)	18 (15%)	4 (3%)	5	24
12	3I	123/132 (93%)	104 (85%)	19 (15%)	0	100	100
13	4A	115/126 (91%)	95 (83%)	19 (16%)	1 (1%)	21	58
13	4I	116/126 (92%)	95 (82%)	20 (17%)	1 (1%)	21	58
14	5A	56/61 (92%)	47 (84%)	8 (14%)	1 (2%)	11	39
14	5I	58/61 (95%)	49 (84%)	7 (12%)	2 (3%)	5	23
15	6A	86/89 (97%)	78 (91%)	8 (9%)	0	100	100
15	6I	86/89 (97%)	77 (90%)	9 (10%)	0	100	100
16	7A	82/88 (93%)	78 (95%)	4 (5%)	0	100	100
16	7I	82/88 (93%)	78 (95%)	4 (5%)	0	100	100
17	8A	98/105 (93%)	89 (91%)	9 (9%)	0	100	100
17	8I	98/105 (93%)	95 (97%)	3 (3%)	0	100	100
18	9A	70/88 (80%)	57 (81%)	12 (17%)	1 (1%)	14	46
18	9I	70/88 (80%)	62 (89%)	6 (9%)	2 (3%)	6	27
19	AA	76/93 (82%)	60 (79%)	14 (18%)	2 (3%)	7	29
19	AI	79/93 (85%)	68 (86%)	8 (10%)	3 (4%)	4	21
20	BA	97/106 (92%)	84 (87%)	12 (12%)	1 (1%)	19	56
20	BI	97/106 (92%)	84 (87%)	12 (12%)	1 (1%)	19	56
21	1B	23/27 (85%)	21 (91%)	2 (9%)	0	100	100
21	1F	23/27 (85%)	22 (96%)	1 (4%)	0	100	100
28	11	270/276 (98%)	247 (92%)	17 (6%)	6 (2%)	8	34
28	19	271/276 (98%)	252 (93%)	16 (6%)	3 (1%)	17	53
29	21	203/206 (98%)	166 (82%)	29 (14%)	8 (4%)	4	20
29	29	203/206 (98%)	160 (79%)	32 (16%)	11 (5%)	2	13
30	31	200/210 (95%)	183 (92%)	16 (8%)	1 (0%)	34	70
30	39	206/210 (98%)	174 (84%)	28 (14%)	4 (2%)	10	38
31	41	179/182 (98%)	155 (87%)	20 (11%)	4 (2%)	8	34
31	49	179/182 (98%)	155 (87%)	22 (12%)	2 (1%)	17	53
32	51	172/180 (96%)	148 (86%)	19 (11%)	5 (3%)	6	27
32	59	167/180 (93%)	136 (81%)	28 (17%)	3 (2%)	11	39

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
33	61	144/148 (97%)	118 (82%)	21 (15%)	5 (4%)	4	23
33	69	144/148 (97%)	116 (81%)	25 (17%)	3 (2%)	9	35
34	15	136/140 (97%)	122 (90%)	14 (10%)	0	100	100
34	58	136/140 (97%)	114 (84%)	19 (14%)	3 (2%)	8	34
35	25	120/122 (98%)	113 (94%)	7 (6%)	0	100	100
35	68	120/122 (98%)	112 (93%)	8 (7%)	0	100	100
36	35	148/150 (99%)	110 (74%)	35 (24%)	3 (2%)	9	36
36	78	148/150 (99%)	115 (78%)	25 (17%)	8 (5%)	2	13
37	45	139/141 (99%)	113 (81%)	22 (16%)	4 (3%)	6	27
37	88	134/141 (95%)	112 (84%)	18 (13%)	4 (3%)	5	26
38	55	115/118 (98%)	108 (94%)	7 (6%)	0	100	100
38	98	116/118 (98%)	100 (86%)	15 (13%)	1 (1%)	21	58
39	65	109/112 (97%)	87 (80%)	19 (17%)	3 (3%)	6	28
39	A8	109/112 (97%)	89 (82%)	19 (17%)	1 (1%)	21	58
40	75	135/146 (92%)	113 (84%)	21 (16%)	1 (1%)	26	64
40	B8	135/146 (92%)	122 (90%)	13 (10%)	0	100	100
41	85	115/118 (98%)	107 (93%)	7 (6%)	1 (1%)	21	58
41	C8	115/118 (98%)	108 (94%)	6 (5%)	1 (1%)	21	58
42	95	99/101 (98%)	81 (82%)	15 (15%)	3 (3%)	5	26
42	D8	99/101 (98%)	90 (91%)	8 (8%)	1 (1%)	19	56
43	A5	111/113 (98%)	103 (93%)	8 (7%)	0	100	100
43	E8	111/113 (98%)	103 (93%)	8 (7%)	0	100	100
44	B5	90/96 (94%)	82 (91%)	7 (8%)	1 (1%)	17	53
44	F8	92/96 (96%)	83 (90%)	6 (6%)	3 (3%)	5	24
45	C5	102/110 (93%)	73 (72%)	24 (24%)	5 (5%)	3	15
45	G8	102/110 (93%)	84 (82%)	14 (14%)	4 (4%)	4	20
46	D5	177/206 (86%)	136 (77%)	31 (18%)	10 (6%)	2	12
46	H8	173/206 (84%)	142 (82%)	27 (16%)	4 (2%)	8	32
47	E5	74/85 (87%)	68 (92%)	6 (8%)	0	100	100
47	I8	78/85 (92%)	66 (85%)	11 (14%)	1 (1%)	15	48
48	F5	95/98 (97%)	83 (87%)	11 (12%)	1 (1%)	17	53

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
48	J8	95/98 (97%)	88 (93%)	7 (7%)	0	100	100
49	G5	64/72 (89%)	57 (89%)	5 (8%)	2 (3%)	5	25
49	K8	65/72 (90%)	56 (86%)	7 (11%)	2 (3%)	5	25
50	H5	57/60 (95%)	53 (93%)	4 (7%)	0	100	100
50	L8	55/60 (92%)	49 (89%)	5 (9%)	1 (2%)	11	39
51	I5	61/71 (86%)	32 (52%)	27 (44%)	2 (3%)	5	24
51	M8	64/71 (90%)	45 (70%)	16 (25%)	3 (5%)	3	16
52	J5	54/60 (90%)	49 (91%)	5 (9%)	0	100	100
52	N8	56/60 (93%)	52 (93%)	4 (7%)	0	100	100
53	K5	43/54 (80%)	29 (67%)	14 (33%)	0	100	100
53	O8	43/54 (80%)	31 (72%)	12 (28%)	0	100	100
54	L5	44/49 (90%)	43 (98%)	1 (2%)	0	100	100
54	P8	43/49 (88%)	41 (95%)	2 (5%)	0	100	100
55	M5	58/65 (89%)	46 (79%)	11 (19%)	1 (2%)	11	41
55	Q8	59/65 (91%)	39 (66%)	16 (27%)	4 (7%)	1	8
All	All	11325/12054 (94%)	9841 (87%)	1312 (12%)	172 (2%)	13	44

All (172) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
28	11	239	ARG
31	41	14	GLU
44	F8	68	ARG
49	K8	47	ASN
49	K8	48	HIS
55	Q8	49	VAL
29	29	25	VAL
30	39	124	LEU
45	C5	29	GLU
46	D5	53	ILE
46	D5	165	VAL
46	D5	171	ILE
48	F5	30	VAL
51	I5	5	ILE
8	7E	86	ILE
18	9I	22	VAL
29	21	78	LEU

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Mol	Chain	Res	Type
32	51	92	ILE
36	78	107	LYS
37	88	66	ILE
37	88	79	LEU
41	C8	93	LYS
46	H8	165	VAL
55	Q8	27	THR
55	Q8	44	LYS
55	Q8	55	ALA
2	12	7	VAL
2	12	71	VAL
12	3A	18	VAL
12	3A	26	ALA
19	AA	11	VAL
28	19	33	LEU
29	29	81	ILE
30	39	84	VAL
30	39	123	LEU
36	35	15	ARG
37	45	135	ASP
39	65	89	ARG
46	D5	105	VAL
4	3E	31	CYS
13	4I	83	ASP
28	11	272	ALA
29	21	60	ASN
31	41	96	ARG
33	61	116	LEU
33	61	145	VAL
34	58	97	ARG
34	58	128	HIS
36	78	6	LEU
38	98	11	ASN
46	H8	6	LYS
11	2A	101	SER
28	19	237	GLU
29	29	9	VAL
29	29	24	THR
29	29	51	PHE
30	39	132	VAL
36	35	6	LEU
36	35	35	HIS

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Mol	Chain	Res	Type
39	65	87	PHE
49	G5	47	ASN
4	3E	81	GLU
4	3E	155	LEU
28	11	237	GLU
28	11	240	ALA
29	21	56	PRO
29	21	118	LYS
31	41	97	ASP
36	78	19	VAL
36	78	23	PRO
39	A8	4	LEU
44	F8	40	LYS
45	G8	81	LYS
45	G8	84	ARG
51	M8	34	GLU
5	42	85	GLY
14	5A	29	ARG
29	29	26	ILE
29	29	90	THR
31	49	47	LYS
32	59	92	ILE
33	69	111	PRO
33	69	145	VAL
37	45	18	LYS
39	65	55	ALA
41	85	93	LYS
42	95	85	LYS
45	C5	17	SER
46	D5	116	VAL
46	D5	161	VAL
49	G5	48	HIS
55	M5	31	HIS
2	1E	156	LYS
2	1E	168	THR
3	2E	15	THR
3	2E	107	GLN
31	41	5	VAL
32	51	10	PRO
32	51	83	TYR
33	61	83	ALA
34	58	22	THR

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Mol	Chain	Res	Type
36	78	16	ARG
37	88	134	ARG
45	G8	83	THR
47	I8	83	PRO
50	L8	54	VAL
29	29	62	PRO
37	45	78	PRO
46	D5	93	ASP
2	1E	135	GLN
3	2E	4	LYS
5	4E	115	VAL
14	5I	13	THR
14	5I	14	PRO
32	51	12	PRO
33	61	12	LEU
33	61	133	HIS
36	78	15	ARG
20	BA	71	THR
40	75	94	ALA
42	95	72	VAL
2	1E	230	VAL
11	2I	82	VAL
19	AI	41	VAL
29	21	55	ASN
44	F8	67	GLY
29	29	91	VAL
37	45	27	VAL
51	I5	33	VAL
20	BI	63	ILE
29	21	72	VAL
36	78	95	VAL
28	19	3	VAL
31	49	5	VAL
44	B5	51	VAL
45	C5	76	CYS
45	C5	85	VAL
2	1E	229	VAL
18	9I	39	VAL
28	11	3	VAL
30	31	132	VAL
46	H8	53	ILE
46	H8	141	VAL

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Mol	Chain	Res	Type
51	M8	50	VAL
2	12	39	ILE
12	3A	47	LYS
13	4A	84	ILE
29	29	59	VAL
2	1E	239	VAL
19	AI	9	VAL
19	AI	67	VAL
29	21	4	ILE
32	51	127	GLU
36	78	7	ARG
37	88	27	VAL
51	M8	5	ILE
12	3A	96	VAL
19	AA	67	VAL
29	29	56	PRO
33	69	144	VAL
42	95	99	ILE
46	D5	61	LEU
46	D5	141	VAL
28	11	123	ALA
42	D8	49	THR
45	G8	76	CYS
18	9A	22	VAL
32	59	8	PRO
32	59	167	GLU
45	C5	3	VAL
46	D5	176	PRO
29	21	147	PRO

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
2	12	205/220 (93%)	167 (82%)	38 (18%)	<div>28</div>
2	1E	205/220 (93%)	167 (82%)	38 (18%)	<div>28</div>

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
3	22	160/188 (85%)	132 (82%)	28 (18%)	2	9
3	2E	159/188 (85%)	131 (82%)	28 (18%)	2	9
4	32	180/181 (99%)	148 (82%)	32 (18%)	2	9
4	3E	180/181 (99%)	147 (82%)	33 (18%)	2	8
5	42	116/123 (94%)	93 (80%)	23 (20%)	1	6
5	4E	116/123 (94%)	87 (75%)	29 (25%)	1	2
6	52	90/90 (100%)	76 (84%)	14 (16%)	3	13
6	5E	90/90 (100%)	78 (87%)	12 (13%)	5	19
7	62	126/127 (99%)	105 (83%)	21 (17%)	3	10
7	6E	126/127 (99%)	103 (82%)	23 (18%)	2	8
8	72	119/119 (100%)	99 (83%)	20 (17%)	2	10
8	7E	119/119 (100%)	97 (82%)	22 (18%)	2	8
9	82	98/99 (99%)	81 (83%)	17 (17%)	2	10
9	8E	98/99 (99%)	76 (78%)	22 (22%)	1	4
10	1A	89/92 (97%)	81 (91%)	8 (9%)	12	39
10	1I	89/92 (97%)	78 (88%)	11 (12%)	6	22
11	2A	88/99 (89%)	76 (86%)	12 (14%)	5	18
11	2I	88/99 (89%)	75 (85%)	13 (15%)	4	15
12	3A	104/109 (95%)	81 (78%)	23 (22%)	1	4
12	3I	104/109 (95%)	88 (85%)	16 (15%)	3	13
13	4A	94/101 (93%)	79 (84%)	15 (16%)	3	12
13	4I	94/101 (93%)	75 (80%)	19 (20%)	1	6
14	5A	48/50 (96%)	41 (85%)	7 (15%)	4	15
14	5I	49/50 (98%)	37 (76%)	12 (24%)	1	2
15	6A	79/80 (99%)	70 (89%)	9 (11%)	7	26
15	6I	79/80 (99%)	71 (90%)	8 (10%)	9	32
16	7A	72/74 (97%)	58 (81%)	14 (19%)	2	7
16	7I	72/74 (97%)	56 (78%)	16 (22%)	1	4
17	8A	95/97 (98%)	83 (87%)	12 (13%)	5	21
17	8I	95/97 (98%)	81 (85%)	14 (15%)	4	15
18	9A	63/77 (82%)	49 (78%)	14 (22%)	1	4

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
18	9I	63/77 (82%)	53 (84%)	10 (16%)	3	12
19	AA	67/80 (84%)	52 (78%)	15 (22%)	1	4
19	AI	70/80 (88%)	49 (70%)	21 (30%)	0	1
20	BA	76/82 (93%)	61 (80%)	15 (20%)	1	7
20	BI	76/82 (93%)	62 (82%)	14 (18%)	2	8
21	1B	20/22 (91%)	18 (90%)	2 (10%)	9	32
21	1F	20/22 (91%)	19 (95%)	1 (5%)	30	66
28	11	214/218 (98%)	171 (80%)	43 (20%)	1	6
28	19	214/218 (98%)	180 (84%)	34 (16%)	3	12
29	21	165/166 (99%)	122 (74%)	43 (26%)	0	1
29	29	165/166 (99%)	134 (81%)	31 (19%)	2	7
30	31	161/166 (97%)	130 (81%)	31 (19%)	2	7
30	39	165/166 (99%)	132 (80%)	33 (20%)	1	6
31	41	155/156 (99%)	128 (83%)	27 (17%)	2	10
31	49	155/156 (99%)	132 (85%)	23 (15%)	4	15
32	51	145/148 (98%)	114 (79%)	31 (21%)	1	5
32	59	141/148 (95%)	112 (79%)	29 (21%)	1	6
33	61	122/124 (98%)	88 (72%)	34 (28%)	0	1
33	69	122/124 (98%)	93 (76%)	29 (24%)	1	3
34	15	117/119 (98%)	95 (81%)	22 (19%)	2	7
34	58	117/119 (98%)	94 (80%)	23 (20%)	1	7
35	25	100/100 (100%)	77 (77%)	23 (23%)	1	4
35	68	100/100 (100%)	84 (84%)	16 (16%)	3	12
36	35	116/116 (100%)	76 (66%)	40 (34%)	0	0
36	78	116/116 (100%)	80 (69%)	36 (31%)	0	0
37	45	111/111 (100%)	84 (76%)	27 (24%)	1	2
37	88	104/111 (94%)	81 (78%)	23 (22%)	1	4
38	55	100/101 (99%)	78 (78%)	22 (22%)	1	4
38	98	101/101 (100%)	81 (80%)	20 (20%)	1	6
39	65	87/88 (99%)	64 (74%)	23 (26%)	0	1
39	A8	87/88 (99%)	60 (69%)	27 (31%)	0	0

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
40	75	120/127 (94%)	92 (77%)	28 (23%)	1	3
40	B8	120/127 (94%)	89 (74%)	31 (26%)	0	2
41	85	93/94 (99%)	80 (86%)	13 (14%)	4	17
41	C8	93/94 (99%)	76 (82%)	17 (18%)	2	8
42	95	82/82 (100%)	57 (70%)	25 (30%)	0	1
42	D8	82/82 (100%)	61 (74%)	21 (26%)	0	2
43	A5	92/92 (100%)	78 (85%)	14 (15%)	3	14
43	E8	92/92 (100%)	72 (78%)	20 (22%)	1	5
44	B5	74/78 (95%)	57 (77%)	17 (23%)	1	4
44	F8	76/78 (97%)	59 (78%)	17 (22%)	1	4
45	C5	85/91 (93%)	61 (72%)	24 (28%)	0	1
45	G8	85/91 (93%)	67 (79%)	18 (21%)	1	5
46	D5	158/179 (88%)	131 (83%)	27 (17%)	2	10
46	H8	154/179 (86%)	123 (80%)	31 (20%)	1	6
47	E5	61/67 (91%)	52 (85%)	9 (15%)	4	15
47	I8	61/67 (91%)	52 (85%)	9 (15%)	4	15
48	F5	82/83 (99%)	67 (82%)	15 (18%)	2	8
48	J8	82/83 (99%)	65 (79%)	17 (21%)	1	6
49	G5	62/67 (92%)	53 (86%)	9 (14%)	4	16
49	K8	62/67 (92%)	41 (66%)	21 (34%)	0	0
50	H5	51/52 (98%)	42 (82%)	9 (18%)	2	9
50	L8	49/52 (94%)	40 (82%)	9 (18%)	2	8
51	I5	57/63 (90%)	47 (82%)	10 (18%)	2	9
51	M8	59/63 (94%)	44 (75%)	15 (25%)	1	2
52	J5	48/52 (92%)	37 (77%)	11 (23%)	1	4
52	N8	51/52 (98%)	38 (74%)	13 (26%)	1	2
53	K5	44/52 (85%)	32 (73%)	12 (27%)	0	1
53	O8	44/52 (85%)	26 (59%)	18 (41%)	0	0
54	L5	39/42 (93%)	32 (82%)	7 (18%)	2	9
54	P8	38/42 (90%)	32 (84%)	6 (16%)	3	12
55	M5	49/55 (89%)	37 (76%)	12 (24%)	1	2

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
55	Q8	50/55 (91%)	32 (64%)	18 (36%)	0	0
All	All	9556/9998 (96%)	7642 (80%)	1914 (20%)	1	6

All (1914) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	1E	6	THR
2	1E	8	LYS
2	1E	9	GLU
2	1E	15	VAL
2	1E	17	PHE
2	1E	24	TRP
2	1E	28	PHE
2	1E	55	PHE
2	1E	63	MET
2	1E	71	VAL
2	1E	74	LYS
2	1E	81	VAL
2	1E	96	ARG
2	1E	107	THR
2	1E	111	ARG
2	1E	113	HIS
2	1E	122	PHE
2	1E	130	ARG
2	1E	136	VAL
2	1E	145	LEU
2	1E	155	LEU
2	1E	157	ARG
2	1E	160	ASP
2	1E	162	ILE
2	1E	164	VAL
2	1E	170	GLU
2	1E	172	ILE
2	1E	178	ARG
2	1E	190	THR
2	1E	196	LEU
2	1E	200	ILE
2	1E	209	ARG
2	1E	213	LEU
2	1E	215	LEU
2	1E	221	LEU
2	1E	226	ARG

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Mol	Chain	Res	Type
2	1E	229	VAL
2	1E	233	SER
3	2E	3	ASN
3	2E	4	LYS
3	2E	5	ILE
3	2E	8	ILE
3	2E	15	THR
3	2E	21	ARG
3	2E	29	TYR
3	2E	34	LEU
3	2E	36	ASP
3	2E	40	ARG
3	2E	45	LYS
3	2E	52	LEU
3	2E	56	ASP
3	2E	62	ASP
3	2E	79	ARG
3	2E	95	THR
3	2E	98	ASN
3	2E	102	ASN
3	2E	127	ARG
3	2E	138	VAL
3	2E	165	THR
3	2E	167	TRP
3	2E	179	ARG
3	2E	190	ARG
3	2E	192	THR
3	2E	196	LEU
3	2E	202	ILE
3	2E	206	GLU
4	3E	3	ARG
4	3E	10	ARG
4	3E	15	GLU
4	3E	24	GLU
4	3E	31	CYS
4	3E	46	LYS
4	3E	47	ARG
4	3E	50	ARG
4	3E	52	SER
4	3E	53	ASP
4	3E	58	LEU
4	3E	66	ARG

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Mol	Chain	Res	Type
4	3E	85	LYS
4	3E	86	LYS
4	3E	89	THR
4	3E	99	SER
4	3E	104	VAL
4	3E	106	TYR
4	3E	108	LEU
4	3E	122	ARG
4	3E	127	THR
4	3E	135	LEU
4	3E	138	TYR
4	3E	141	ARG
4	3E	156	GLU
4	3E	175	SER
4	3E	179	GLU
4	3E	184	LYS
4	3E	190	ASP
4	3E	191	ARG
4	3E	193	ASP
4	3E	200	GLU
4	3E	209	ARG
5	4E	5	ASP
5	4E	6	PHE
5	4E	10	MET
5	4E	11	ILE
5	4E	13	ILE
5	4E	14	ARG
5	4E	16	THR
5	4E	18	ARG
5	4E	33	VAL
5	4E	41	VAL
5	4E	50	GLU
5	4E	56	GLN
5	4E	63	ARG
5	4E	64	ARG
5	4E	68	GLU
5	4E	72	GLN
5	4E	73	ASN
5	4E	75	THR
5	4E	79	GLU
5	4E	81	GLU
5	4E	87	SER

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Mol	Chain	Res	Type
5	4E	91	LEU
5	4E	112	LEU
5	4E	116	THR
5	4E	126	ARG
5	4E	131	ILE
5	4E	151	LEU
5	4E	152	ARG
5	4E	153	LYS
6	5E	17	SER
6	5E	23	LYS
6	5E	27	GLN
6	5E	41	GLU
6	5E	46	ARG
6	5E	55	ASP
6	5E	64	GLN
6	5E	65	VAL
6	5E	70	ASP
6	5E	75	LEU
6	5E	89	MET
6	5E	91	VAL
7	6E	6	ARG
7	6E	8	GLU
7	6E	10	ARG
7	6E	36	LYS
7	6E	38	LEU
7	6E	45	ASP
7	6E	47	CYS
7	6E	48	LYS
7	6E	54	THR
7	6E	59	LEU
7	6E	63	LYS
7	6E	66	VAL
7	6E	79	ARG
7	6E	80	VAL
7	6E	85	TYR
7	6E	90	GLU
7	6E	91	VAL
7	6E	104	LEU
7	6E	113	GLU
7	6E	115	ARG
7	6E	124	LEU
7	6E	155	ARG

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Mol	Chain	Res	Type
7	6E	156	TRP
8	7E	1	MET
8	7E	29	SER
8	7E	30	ARG
8	7E	37	ARG
8	7E	45	ILE
8	7E	52	ASP
8	7E	54	ASP
8	7E	68	ARG
8	7E	80	ILE
8	7E	82	HIS
8	7E	83	ILE
8	7E	85	ARG
8	7E	88	LYS
8	7E	91	ARG
8	7E	95	VAL
8	7E	102	ARG
8	7E	104	ARG
8	7E	107	LEU
8	7E	112	LEU
8	7E	129	VAL
8	7E	134	ILE
8	7E	137	VAL
9	8E	9	ARG
9	8E	10	ARG
9	8E	20	ARG
9	8E	38	GLN
9	8E	41	VAL
9	8E	42	ARG
9	8E	47	LEU
9	8E	53	VAL
9	8E	64	THR
9	8E	79	LEU
9	8E	89	ASN
9	8E	91	ASP
9	8E	92	TYR
9	8E	95	LYS
9	8E	104	ARG
9	8E	105	ASP
9	8E	108	VAL
9	8E	112	LYS
9	8E	117	HIS

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Mol	Chain	Res	Type
9	8E	118	LYS
9	8E	121	ARG
9	8E	125	TYR
10	1I	5	ARG
10	1I	54	PHE
10	1I	58	ASP
10	1I	62	HIS
10	1I	66	ARG
10	1I	70	ARG
10	1I	76	ASN
10	1I	92	THR
10	1I	96	ILE
10	1I	98	ILE
10	1I	101	VAL
11	2I	14	VAL
11	2I	31	THR
11	2I	32	ILE
11	2I	36	ASP
11	2I	81	ASP
11	2I	87	THR
11	2I	91	ARG
11	2I	103	LEU
11	2I	105	VAL
11	2I	106	LYS
11	2I	109	VAL
11	2I	114	VAL
11	2I	116	HIS
12	3I	11	VAL
12	3I	33	ARG
12	3I	34	ARG
12	3I	44	THR
12	3I	50	SER
12	3I	60	LEU
12	3I	62	SER
12	3I	65	GLU
12	3I	67	THR
12	3I	85	ILE
12	3I	89	ARG
12	3I	91	LYS
12	3I	100	ILE
12	3I	102	ARG
12	3I	114	LYS

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Mol	Chain	Res	Type
12	3I	116	SER
13	4I	3	ARG
13	4I	9	ILE
13	4I	11	ARG
13	4I	13	LYS
13	4I	19	LEU
13	4I	34	LEU
13	4I	48	LEU
13	4I	56	LEU
13	4I	64	TRP
13	4I	70	LEU
13	4I	77	ASN
13	4I	98	VAL
13	4I	99	ARG
13	4I	105	THR
13	4I	106	ASN
13	4I	108	ARG
13	4I	110	ARG
13	4I	111	LYS
13	4I	117	VAL
14	5I	12	ARG
14	5I	17	LYS
14	5I	18	VAL
14	5I	22	THR
14	5I	23	ARG
14	5I	26	ARG
14	5I	32	SER
14	5I	33	VAL
14	5I	41	ARG
14	5I	44	LEU
14	5I	58	LYS
14	5I	60	SER
15	6I	6	GLU
15	6I	22	THR
15	6I	38	ARG
15	6I	41	GLU
15	6I	47	LYS
15	6I	66	LEU
15	6I	67	LEU
15	6I	82	ILE
16	7I	2	VAL
16	7I	6	LEU

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Mol	Chain	Res	Type
16	7I	8	ARG
16	7I	20	VAL
16	7I	27	LYS
16	7I	28	ARG
16	7I	33	ILE
16	7I	35	LYS
16	7I	36	ILE
16	7I	45	THR
16	7I	50	LYS
16	7I	54	GLU
16	7I	67	THR
16	7I	69	THR
16	7I	72	ARG
16	7I	83	GLU
17	8I	38	ARG
17	8I	45	HIS
17	8I	48	GLU
17	8I	50	LYS
17	8I	52	LYS
17	8I	60	ILE
17	8I	62	SER
17	8I	68	ARG
17	8I	74	LEU
17	8I	85	VAL
17	8I	89	LEU
17	8I	92	ARG
17	8I	99	SER
17	8I	101	ARG
18	9I	19	LYS
18	9I	23	LYS
18	9I	25	THR
18	9I	26	LEU
18	9I	31	LEU
18	9I	32	ARG
18	9I	54	ARG
18	9I	56	THR
18	9I	59	SER
18	9I	82	THR
19	AI	4	SER
19	AI	5	LEU
19	AI	7	LYS
19	AI	14	HIS

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Mol	Chain	Res	Type
19	AI	20	LEU
19	AI	21	GLU
19	AI	22	LEU
19	AI	29	ARG
19	AI	30	LEU
19	AI	31	ILE
19	AI	32	LYS
19	AI	37	ARG
19	AI	43	GLU
19	AI	60	VAL
19	AI	61	TYR
19	AI	63	THR
19	AI	65	ASN
19	AI	67	VAL
19	AI	71	LEU
19	AI	77	THR
19	AI	78	ARG
20	BI	9	ASN
20	BI	10	LEU
20	BI	11	SER
20	BI	13	LEU
20	BI	15	ARG
20	BI	37	SER
20	BI	51	GLU
20	BI	57	ARG
20	BI	64	ASP
20	BI	68	LYS
20	BI	73	HIS
20	BI	82	SER
20	BI	99	LEU
20	BI	105	SER
21	1F	6	ARG
28	11	13	ARG
28	11	15	PHE
28	11	17	THR
28	11	31	LYS
28	11	37	LEU
28	11	46	GLN
28	11	58	HIS
28	11	61	LEU
28	11	64	ILE
28	11	65	ILE

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Mol	Chain	Res	Type
28	11	71	ASP
28	11	89	SER
28	11	94	LEU
28	11	95	LEU
28	11	99	ASP
28	11	103	ARG
28	11	105	ILE
28	11	106	ILE
28	11	111	LEU
28	11	115	GLN
28	11	126	GLN
28	11	136	ILE
28	11	142	VAL
28	11	147	LEU
28	11	155	LEU
28	11	157	ARG
28	11	162	SER
28	11	165	ILE
28	11	183	ARG
28	11	192	THR
28	11	193	VAL
28	11	200	ASP
28	11	212	SER
28	11	217	ARG
28	11	221	VAL
28	11	229	VAL
28	11	242	ARG
28	11	257	LEU
28	11	259	THR
28	11	260	ARG
28	11	263	ARG
28	11	271	ILE
28	11	273	ARG
29	21	5	LEU
29	21	13	ARG
29	21	14	ILE
29	21	16	ARG
29	21	17	ASP
29	21	25	VAL
29	21	26	ILE
29	21	38	THR
29	21	40	GLU

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Mol	Chain	Res	Type
29	21	41	LYS
29	21	45	THR
29	21	47	VAL
29	21	48	GLN
29	21	52	LEU
29	21	54	GLN
29	21	55	ASN
29	21	59	VAL
29	21	63	LEU
29	21	67	PHE
29	21	69	LYS
29	21	72	VAL
29	21	78	LEU
29	21	79	ARG
29	21	92	THR
29	21	101	ARG
29	21	107	THR
29	21	111	ARG
29	21	116	VAL
29	21	118	LYS
29	21	119	ARG
29	21	128	SER
29	21	144	ARG
29	21	146	THR
29	21	163	GLU
29	21	167	VAL
29	21	175	VAL
29	21	179	GLU
29	21	185	LYS
29	21	188	VAL
29	21	195	LEU
29	21	197	ILE
29	21	202	LYS
29	21	203	LYS
30	31	8	GLN
30	31	9	ILE
30	31	15	SER
30	31	17	ARG
30	31	18	ARG
30	31	33	LEU
30	31	36	VAL
30	31	38	ARG

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Mol	Chain	Res	Type
30	31	64	ILE
30	31	67	GLN
30	31	74	ARG
30	31	78	ILE
30	31	88	VAL
30	31	98	SER
30	31	106	ARG
30	31	116	ASP
30	31	117	ARG
30	31	127	GLU
30	31	140	LEU
30	31	153	SER
30	31	158	THR
30	31	164	ARG
30	31	165	ARG
30	31	176	LEU
30	31	181	LEU
30	31	183	VAL
30	31	188	ARG
30	31	191	ARG
30	31	192	LEU
30	31	201	VAL
30	31	204	ASN
31	41	3	LEU
31	41	10	LYS
31	41	19	LEU
31	41	20	ILE
31	41	28	VAL
31	41	31	VAL
31	41	43	LEU
31	41	45	GLU
31	41	54	GLU
31	41	58	GLN
31	41	62	LEU
31	41	63	ILE
31	41	67	LYS
31	41	76	SER
31	41	80	PHE
31	41	82	LEU
31	41	86	MET
31	41	90	LEU
31	41	94	LEU

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Mol	Chain	Res	Type
31	41	101	ILE
31	41	116	ASP
31	41	128	ARG
31	41	133	LEU
31	41	155	MET
31	41	156	ASP
31	41	165	THR
31	41	166	ASP
32	51	3	ARG
32	51	4	ILE
32	51	7	LEU
32	51	10	PRO
32	51	24	VAL
32	51	37	VAL
32	51	40	GLU
32	51	41	MET
32	51	45	VAL
32	51	50	VAL
32	51	64	LEU
32	51	71	LEU
32	51	72	ILE
32	51	77	LYS
32	51	80	SER
32	51	81	GLU
32	51	83	TYR
32	51	88	LEU
32	51	92	ILE
32	51	95	ARG
32	51	104	GLU
32	51	106	THR
32	51	122	THR
32	51	129	THR
32	51	132	ARG
32	51	136	ILE
32	51	139	GLN
32	51	141	VAL
32	51	149	ARG
32	51	153	LYS
32	51	170	ARG
33	61	2	LYS
33	61	9	LEU
33	61	20	ASP

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Mol	Chain	Res	Type
33	61	38	LEU
33	61	40	THR
33	61	41	GLU
33	61	42	SER
33	61	48	GLU
33	61	56	LYS
33	61	58	LEU
33	61	60	GLU
33	61	64	GLU
33	61	74	ASN
33	61	77	LEU
33	61	78	THR
33	61	79	ILE
33	61	81	VAL
33	61	82	ARG
33	61	85	GLU
33	61	86	THR
33	61	88	ILE
33	61	92	VAL
33	61	95	LYS
33	61	105	HIS
33	61	107	VAL
33	61	110	ASP
33	61	117	GLU
33	61	122	GLU
33	61	131	LYS
33	61	135	GLU
33	61	140	LEU
33	61	142	VAL
33	61	144	VAL
33	61	145	VAL
34	58	2	LYS
34	58	5	VAL
34	58	7	LYS
34	58	10	GLU
34	58	12	ARG
34	58	28	THR
34	58	34	LEU
34	58	38	HIS
34	58	43	THR
34	58	48	MET
34	58	58	ASP

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Mol	Chain	Res	Type
34	58	60	ILE
34	58	61	ARG
34	58	65	LYS
34	58	67	LEU
34	58	90	MET
34	58	97	ARG
34	58	99	LEU
34	58	120	LEU
34	58	128	HIS
34	58	133	GLN
34	58	134	ARG
34	58	136	GLU
35	68	23	ARG
35	68	24	VAL
35	68	28	SER
35	68	32	TYR
35	68	38	VAL
35	68	53	LYS
35	68	68	GLU
35	68	75	SER
35	68	80	ASP
35	68	91	LEU
35	68	94	ARG
35	68	98	VAL
35	68	112	MET
35	68	113	LYS
35	68	115	VAL
35	68	116	SER
36	78	2	LYS
36	78	4	SER
36	78	6	LEU
36	78	7	ARG
36	78	10	PRO
36	78	13	ASN
36	78	15	ARG
36	78	18	ARG
36	78	19	VAL
36	78	29	LYS
36	78	30	THR
36	78	32	THR
36	78	39	LYS
36	78	41	ARG

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Mol	Chain	Res	Type
36	78	45	LEU
36	78	46	LYS
36	78	49	ARG
36	78	50	ARG
36	78	57	THR
36	78	75	ILE
36	78	76	LYS
36	78	77	ARG
36	78	88	LEU
36	78	100	LEU
36	78	105	LEU
36	78	106	LEU
36	78	108	LYS
36	78	112	LEU
36	78	123	LEU
36	78	135	LEU
36	78	138	LEU
36	78	144	GLU
36	78	146	VAL
36	78	147	LEU
36	78	148	LEU
36	78	149	GLU
37	88	1	MET
37	88	5	ARG
37	88	6	ARG
37	88	18	LYS
37	88	25	ASP
37	88	35	VAL
37	88	45	GLN
37	88	51	ARG
37	88	55	VAL
37	88	56	ARG
37	88	58	PHE
37	88	59	ARG
37	88	60	ARG
37	88	64	ILE
37	88	67	ARG
37	88	82	ARG
37	88	109	VAL
37	88	110	THR
37	88	112	GLU
37	88	119	ARG

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Mol	Chain	Res	Type
37	88	134	ARG
37	88	139	GLU
37	88	141	GLN
38	98	4	LEU
38	98	9	LYS
38	98	10	LEU
38	98	12	ARG
38	98	15	SER
38	98	18	LEU
38	98	28	LEU
38	98	29	LEU
38	98	34	ILE
38	98	35	THR
38	98	44	LEU
38	98	45	ARG
38	98	59	ASP
38	98	65	LEU
38	98	79	LEU
38	98	91	GLN
38	98	95	THR
38	98	96	ARG
38	98	104	ARG
38	98	118	GLU
39	A8	3	ARG
39	A8	4	LEU
39	A8	15	ARG
39	A8	17	ARG
39	A8	19	LYS
39	A8	20	ARG
39	A8	24	LEU
39	A8	29	PHE
39	A8	30	ARG
39	A8	35	ILE
39	A8	36	TYR
39	A8	46	VAL
39	A8	50	SER
39	A8	52	SER
39	A8	54	LEU
39	A8	56	LEU
39	A8	57	LYS
39	A8	61	ASN
39	A8	73	LEU

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Mol	Chain	Res	Type
39	A8	84	GLN
39	A8	89	ARG
39	A8	95	HIS
39	A8	98	VAL
39	A8	101	LEU
39	A8	106	ARG
39	A8	107	GLU
39	A8	112	PHE
40	B8	7	ILE
40	B8	10	VAL
40	B8	12	SER
40	B8	13	ARG
40	B8	16	ARG
40	B8	21	GLU
40	B8	23	ARG
40	B8	27	THR
40	B8	30	VAL
40	B8	33	LYS
40	B8	39	ARG
40	B8	41	ARG
40	B8	42	ILE
40	B8	50	ILE
40	B8	58	ASN
40	B8	59	THR
40	B8	62	THR
40	B8	64	ARG
40	B8	65	LYS
40	B8	74	ARG
40	B8	85	LYS
40	B8	86	ILE
40	B8	89	VAL
40	B8	99	LEU
40	B8	105	LEU
40	B8	106	SER
40	B8	108	ARG
40	B8	110	ILE
40	B8	111	ARG
40	B8	112	ARG
40	B8	128	GLU
41	C8	3	ARG
41	C8	5	LYS
41	C8	11	ARG

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Mol	Chain	Res	Type
41	C8	27	LEU
41	C8	34	LYS
41	C8	52	ARG
41	C8	57	PHE
41	C8	59	ARG
41	C8	69	CYS
41	C8	74	LEU
41	C8	79	PHE
41	C8	89	GLU
41	C8	92	ARG
41	C8	94	ASN
41	C8	95	LEU
41	C8	104	GLN
41	C8	111	GLU
42	D8	1	MET
42	D8	6	LYS
42	D8	7	THR
42	D8	18	LEU
42	D8	20	LEU
42	D8	24	LYS
42	D8	34	GLU
42	D8	35	LEU
42	D8	38	LEU
42	D8	39	LEU
42	D8	40	LEU
42	D8	45	THR
42	D8	47	VAL
42	D8	49	THR
42	D8	58	VAL
42	D8	64	HIS
42	D8	73	SER
42	D8	79	VAL
42	D8	88	ARG
42	D8	91	TYR
42	D8	100	ARG
43	E8	11	ARG
43	E8	12	ILE
43	E8	13	SER
43	E8	20	VAL
43	E8	39	THR
43	E8	51	LEU
43	E8	60	ASN

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Mol	Chain	Res	Type
43	E8	66	GLU
43	E8	69	LEU
43	E8	70	TYR
43	E8	76	VAL
43	E8	78	GLU
43	E8	84	ARG
43	E8	88	ARG
43	E8	90	ARG
43	E8	92	ARG
43	E8	96	ILE
43	E8	103	ILE
43	E8	107	LEU
43	E8	111	HIS
44	F8	2	LYS
44	F8	12	VAL
44	F8	15	GLU
44	F8	28	PHE
44	F8	40	LYS
44	F8	49	VAL
44	F8	53	LYS
44	F8	54	VAL
44	F8	57	LEU
44	F8	65	ARG
44	F8	66	LEU
44	F8	68	ARG
44	F8	70	LEU
44	F8	72	LYS
44	F8	80	ILE
44	F8	81	VAL
44	F8	88	LYS
45	G8	4	LYS
45	G8	6	HIS
45	G8	14	LEU
45	G8	38	ILE
45	G8	44	ILE
45	G8	45	VAL
45	G8	52	SER
45	G8	57	GLN
45	G8	61	ILE
45	G8	67	LEU
45	G8	71	LYS
45	G8	79	CYS

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Mol	Chain	Res	Type
45	G8	82	PRO
45	G8	84	ARG
45	G8	85	VAL
45	G8	86	ARG
45	G8	96	ILE
45	G8	97	ARG
46	H8	2	GLU
46	H8	11	GLU
46	H8	19	ARG
46	H8	34	ASN
46	H8	37	VAL
46	H8	43	GLU
46	H8	49	ARG
46	H8	53	ILE
46	H8	58	VAL
46	H8	59	LEU
46	H8	61	LEU
46	H8	71	VAL
46	H8	76	LEU
46	H8	77	ASP
46	H8	80	ARG
46	H8	81	ARG
46	H8	82	ARG
46	H8	86	VAL
46	H8	94	GLU
46	H8	103	ARG
46	H8	105	VAL
46	H8	112	ARG
46	H8	117	LEU
46	H8	119	GLU
46	H8	121	HIS
46	H8	132	ASN
46	H8	140	ASP
46	H8	144	LEU
46	H8	149	SER
46	H8	154	ASP
46	H8	166	SER
47	I8	11	ARG
47	I8	36	ILE
47	I8	38	VAL
47	I8	41	ARG
47	I8	53	MET

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Mol	Chain	Res	Type
47	I8	67	VAL
47	I8	70	GLN
47	I8	74	ARG
47	I8	80	HIS
48	J8	19	GLN
48	J8	21	ARG
48	J8	26	ARG
48	J8	33	LYS
48	J8	41	ARG
48	J8	52	ARG
48	J8	65	SER
48	J8	74	VAL
48	J8	78	LYS
48	J8	80	LEU
48	J8	81	LYS
48	J8	86	SER
48	J8	90	ILE
48	J8	91	LYS
48	J8	92	LYS
48	J8	93	GLU
48	J8	94	LEU
49	K8	5	GLU
49	K8	9	GLN
49	K8	14	ARG
49	K8	16	LEU
49	K8	17	SER
49	K8	19	VAL
49	K8	24	LEU
49	K8	32	LEU
49	K8	41	ILE
49	K8	46	GLN
49	K8	47	ASN
49	K8	48	HIS
49	K8	50	ILE
49	K8	53	LEU
49	K8	54	LYS
49	K8	55	ARG
49	K8	59	ARG
49	K8	62	THR
49	K8	64	LEU
49	K8	65	ASN
49	K8	66	GLU

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Mol	Chain	Res	Type
50	L8	6	VAL
50	L8	8	LEU
50	L8	11	SER
50	L8	31	LEU
50	L8	32	GLN
50	L8	37	LEU
50	L8	38	GLU
50	L8	40	THR
50	L8	53	LEU
51	M8	1	MET
51	M8	21	VAL
51	M8	27	THR
51	M8	36	CYS
51	M8	38	LYS
51	M8	39	CYS
51	M8	42	PHE
51	M8	43	TYR
51	M8	44	THR
51	M8	48	ARG
51	M8	55	ARG
51	M8	59	PHE
51	M8	61	ARG
51	M8	62	ARG
51	M8	65	ASP
52	N8	3	LYS
52	N8	6	VAL
52	N8	16	ARG
52	N8	26	THR
52	N8	29	THR
52	N8	36	CYS
52	N8	40	LYS
52	N8	48	GLU
52	N8	49	CYS
52	N8	51	TYR
52	N8	55	ARG
52	N8	56	LYS
52	N8	57	VAL
53	O8	10	LEU
53	O8	12	GLU
53	O8	15	GLU
53	O8	16	CYS
53	O8	26	ASN

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Mol	Chain	Res	Type
53	O8	27	LYS
53	O8	28	ARG
53	O8	30	THR
53	O8	32	ASN
53	O8	33	LYS
53	O8	34	LEU
53	O8	37	ARG
53	O8	39	TYR
53	O8	42	TRP
53	O8	44	ARG
53	O8	47	THR
53	O8	51	GLU
53	O8	52	VAL
54	P8	4	THR
54	P8	8	ASN
54	P8	14	LYS
54	P8	22	MET
54	P8	23	ARG
54	P8	43	THR
55	Q8	8	LYS
55	Q8	19	SER
55	Q8	21	LYS
55	Q8	23	VAL
55	Q8	30	ARG
55	Q8	31	HIS
55	Q8	32	LEU
55	Q8	34	TRP
55	Q8	35	GLN
55	Q8	39	LYS
55	Q8	42	ARG
55	Q8	46	ARG
55	Q8	47	LYS
55	Q8	48	PHE
55	Q8	52	LYS
55	Q8	58	ILE
55	Q8	59	LYS
55	Q8	61	LEU
2	12	4	GLU
2	12	5	ILE
2	12	12	GLU
2	12	17	PHE
2	12	23	ARG

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Mol	Chain	Res	Type
2	12	24	TRP
2	12	31	TYR
2	12	42	ILE
2	12	51	LEU
2	12	52	GLU
2	12	55	PHE
2	12	58	ILE
2	12	69	LEU
2	12	71	VAL
2	12	75	LYS
2	12	78	GLN
2	12	83	MET
2	12	90	MET
2	12	103	THR
2	12	108	ILE
2	12	109	SER
2	12	121	LEU
2	12	122	PHE
2	12	130	ARG
2	12	145	LEU
2	12	155	LEU
2	12	160	ASP
2	12	165	VAL
2	12	178	ARG
2	12	179	LYS
2	12	185	ILE
2	12	187	LEU
2	12	191	ASP
2	12	204	ASN
2	12	205	ASP
2	12	209	ARG
2	12	212	GLN
2	12	233	SER
3	22	5	ILE
3	22	16	ARG
3	22	18	TRP
3	22	21	ARG
3	22	22	TRP
3	22	28	GLN
3	22	29	TYR
3	22	40	ARG
3	22	43	LEU

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Mol	Chain	Res	Type
3	22	47	LEU
3	22	48	TYR
3	22	67	THR
3	22	79	ARG
3	22	83	ARG
3	22	94	LEU
3	22	95	THR
3	22	101	LEU
3	22	104	GLN
3	22	105	GLU
3	22	107	GLN
3	22	119	ARG
3	22	128	PHE
3	22	138	VAL
3	22	140	ARG
3	22	190	ARG
3	22	191	THR
3	22	196	LEU
3	22	202	ILE
4	32	4	TYR
4	32	5	ILE
4	32	8	VAL
4	32	13	ARG
4	32	24	GLU
4	32	28	SER
4	32	30	LYS
4	32	36	ARG
4	32	50	ARG
4	32	58	LEU
4	32	61	LYS
4	32	73	ARG
4	32	76	ARG
4	32	78	LEU
4	32	83	SER
4	32	94	LEU
4	32	107	ARG
4	32	119	GLN
4	32	122	ARG
4	32	126	ILE
4	32	127	THR
4	32	135	LEU
4	32	141	ARG

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Mol	Chain	Res	Type
4	32	157	LEU
4	32	158	ILE
4	32	159	ARG
4	32	162	LEU
4	32	178	VAL
4	32	187	ARG
4	32	191	ARG
4	32	194	LEU
4	32	196	LEU
5	42	13	ILE
5	42	16	THR
5	42	24	ARG
5	42	34	VAL
5	42	41	VAL
5	42	43	LEU
5	42	47	LYS
5	42	61	TYR
5	42	72	GLN
5	42	75	THR
5	42	78	HIS
5	42	79	GLU
5	42	83	GLU
5	42	90	VAL
5	42	101	ILE
5	42	115	VAL
5	42	116	THR
5	42	118	ILE
5	42	126	ARG
5	42	136	MET
5	42	137	GLU
5	42	144	THR
5	42	150	ARG
6	52	3	ARG
6	52	7	ASN
6	52	14	LEU
6	52	16	GLN
6	52	21	LEU
6	52	24	GLU
6	52	28	ARG
6	52	40	VAL
6	52	43	LEU
6	52	54	LYS

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Mol	Chain	Res	Type
6	52	64	GLN
6	52	70	ASP
6	52	71	ARG
6	52	74	ASP
7	62	8	GLU
7	62	21	VAL
7	62	29	LYS
7	62	35	LYS
7	62	36	LYS
7	62	41	ARG
7	62	50	ILE
7	62	52	GLU
7	62	61	VAL
7	62	63	LYS
7	62	72	ARG
7	62	78	ARG
7	62	84	ASN
7	62	90	GLU
7	62	91	VAL
7	62	92	SER
7	62	94	ARG
7	62	104	LEU
7	62	114	ARG
7	62	131	LYS
7	62	140	ASP
8	72	1	MET
8	72	2	LEU
8	72	23	SER
8	72	24	THR
8	72	25	ASP
8	72	33	GLU
8	72	52	ASP
8	72	54	ASP
8	72	77	GLU
8	72	87	SER
8	72	91	ARG
8	72	92	ARG
8	72	95	VAL
8	72	97	VAL
8	72	99	GLU
8	72	100	ILE
8	72	104	ARG

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Mol	Chain	Res	Type
8	72	112	LEU
8	72	116	LYS
8	72	121	ASP
9	82	10	ARG
9	82	27	THR
9	82	38	GLN
9	82	42	ARG
9	82	56	LEU
9	82	58	HIS
9	82	88	TYR
9	82	91	ASP
9	82	95	LYS
9	82	104	ARG
9	82	111	ARG
9	82	113	LYS
9	82	114	TYR
9	82	117	HIS
9	82	118	LYS
9	82	125	TYR
9	82	128	ARG
10	1A	13	HIS
10	1A	17	ASP
10	1A	47	PHE
10	1A	59	SER
10	1A	62	HIS
10	1A	70	ARG
10	1A	79	ARG
10	1A	95	GLU
11	2A	29	ILE
11	2A	30	VAL
11	2A	31	THR
11	2A	57	THR
11	2A	63	LEU
11	2A	95	ILE
11	2A	104	GLN
11	2A	105	VAL
11	2A	106	LYS
11	2A	107	SER
11	2A	119	CYS
11	2A	124	LYS
12	3A	6	THR
12	3A	13	LYS

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Mol	Chain	Res	Type
12	3A	24	VAL
12	3A	27	LEU
12	3A	33	ARG
12	3A	34	ARG
12	3A	37	CYS
12	3A	39	VAL
12	3A	41	ARG
12	3A	42	THR
12	3A	44	THR
12	3A	46	LYS
12	3A	54	LYS
12	3A	57	LYS
12	3A	60	LEU
12	3A	62	SER
12	3A	78	GLN
12	3A	83	VAL
12	3A	84	LEU
12	3A	92	ASP
12	3A	102	ARG
12	3A	118	SER
12	3A	123	LYS
13	4A	8	GLU
13	4A	32	GLU
13	4A	37	THR
13	4A	47	ASP
13	4A	55	ARG
13	4A	64	TRP
13	4A	66	LEU
13	4A	77	ASN
13	4A	82	MET
13	4A	83	ASP
13	4A	91	ARG
13	4A	98	VAL
13	4A	101	GLN
13	4A	108	ARG
13	4A	117	VAL
14	5A	8	GLU
14	5A	12	ARG
14	5A	17	LYS
14	5A	22	THR
14	5A	27	CYS
14	5A	33	VAL

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Mol	Chain	Res	Type
14	5A	44	LEU
15	6A	3	ILE
15	6A	22	THR
15	6A	31	LEU
15	6A	34	LEU
15	6A	41	GLU
15	6A	48	LYS
15	6A	68	ARG
15	6A	87	ILE
15	6A	88	ARG
16	7A	2	VAL
16	7A	6	LEU
16	7A	8	ARG
16	7A	19	ILE
16	7A	21	VAL
16	7A	27	LYS
16	7A	33	ILE
16	7A	47	ASP
16	7A	55	ARG
16	7A	61	SER
16	7A	65	GLN
16	7A	67	THR
16	7A	81	ARG
16	7A	82	GLN
17	8A	4	LYS
17	8A	10	VAL
17	8A	13	ASP
17	8A	49	GLU
17	8A	50	LYS
17	8A	60	ILE
17	8A	62	SER
17	8A	63	ARG
17	8A	70	ARG
17	8A	74	LEU
17	8A	100	LYS
17	8A	101	ARG
18	9A	17	SER
18	9A	26	LEU
18	9A	28	GLU
18	9A	29	PHE
18	9A	32	ARG
18	9A	35	ARG

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Mol	Chain	Res	Type
18	9A	36	ASN
18	9A	42	ARG
18	9A	44	LEU
18	9A	47	THR
18	9A	53	ARG
18	9A	54	ARG
18	9A	59	SER
18	9A	84	LYS
19	AA	11	VAL
19	AA	15	LEU
19	AA	23	ASN
19	AA	25	LYS
19	AA	30	LEU
19	AA	33	THR
19	AA	34	TRP
19	AA	35	SER
19	AA	49	ILE
19	AA	53	ASN
19	AA	60	VAL
19	AA	66	MET
19	AA	71	LEU
19	AA	78	ARG
19	AA	83	HIS
20	BA	10	LEU
20	BA	11	SER
20	BA	23	ARG
20	BA	24	LEU
20	BA	26	ASN
20	BA	37	SER
20	BA	56	MET
20	BA	60	GLU
20	BA	62	LEU
20	BA	64	ASP
20	BA	74	LYS
20	BA	75	ASN
20	BA	83	ARG
20	BA	84	LEU
20	BA	88	VAL
21	1B	9	ARG
21	1B	22	ARG
28	19	20	ASP
28	19	24	ILE

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Mol	Chain	Res	Type
28	19	28	GLU
28	19	30	GLU
28	19	34	VAL
28	19	37	LEU
28	19	43	ARG
28	19	49	ILE
28	19	61	LEU
28	19	64	ILE
28	19	65	ILE
28	19	88	ARG
28	19	94	LEU
28	19	105	ILE
28	19	111	LEU
28	19	138	VAL
28	19	141	VAL
28	19	147	LEU
28	19	155	LEU
28	19	173	VAL
28	19	176	ARG
28	19	182	LEU
28	19	208	LYS
28	19	211	ARG
28	19	217	ARG
28	19	239	ARG
28	19	242	ARG
28	19	244	ARG
28	19	255	LYS
28	19	257	LEU
28	19	260	ARG
28	19	262	ARG
28	19	266	SER
28	19	271	ILE
29	29	5	LEU
29	29	25	VAL
29	29	27	LEU
29	29	37	ARG
29	29	38	THR
29	29	44	TYR
29	29	45	THR
29	29	54	GLN
29	29	60	ASN
29	29	61	ARG

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Mol	Chain	Res	Type
29	29	69	LYS
29	29	75	VAL
29	29	77	ILE
29	29	79	ARG
29	29	87	GLU
29	29	90	THR
29	29	93	VAL
29	29	104	VAL
29	29	107	THR
29	29	111	ARG
29	29	116	VAL
29	29	117	MET
29	29	119	ARG
29	29	144	ARG
29	29	154	LYS
29	29	170	LEU
29	29	175	VAL
29	29	178	GLU
29	29	179	GLU
29	29	197	ILE
29	29	201	THR
30	39	2	LYS
30	39	7	TYR
30	39	8	GLN
30	39	11	VAL
30	39	19	GLU
30	39	20	LEU
30	39	24	LEU
30	39	28	ILE
30	39	38	ARG
30	39	57	VAL
30	39	62	ARG
30	39	63	LYS
30	39	64	ILE
30	39	66	PRO
30	39	67	GLN
30	39	68	LYS
30	39	70	THR
30	39	74	ARG
30	39	77	ASP
30	39	82	ILE
30	39	83	PHE

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Mol	Chain	Res	Type
30	39	98	SER
30	39	106	ARG
30	39	112	MET
30	39	140	LEU
30	39	153	SER
30	39	158	THR
30	39	165	ARG
30	39	191	ARG
30	39	193	VAL
30	39	196	LEU
30	39	201	VAL
30	39	205	ARG
31	49	4	ASP
31	49	7	LEU
31	49	18	GLU
31	49	19	LEU
31	49	28	VAL
31	49	33	ARG
31	49	64	THR
31	49	67	LYS
31	49	71	THR
31	49	74	LYS
31	49	77	ILE
31	49	80	PHE
31	49	82	LEU
31	49	91	ARG
31	49	115	ARG
31	49	118	ARG
31	49	126	ASP
31	49	130	ASN
31	49	133	LEU
31	49	139	LEU
31	49	153	ARG
31	49	157	ILE
31	49	161	THR
32	59	4	ILE
32	59	6	ARG
32	59	7	LEU
32	59	11	VAL
32	59	24	VAL
32	59	27	LYS
32	59	32	GLU

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Mol	Chain	Res	Type
32	59	41	MET
32	59	43	VAL
32	59	56	SER
32	59	70	THR
32	59	85	LYS
32	59	86	GLU
32	59	89	ILE
32	59	101	ARG
32	59	103	LEU
32	59	105	LEU
32	59	107	VAL
32	59	119	GLU
32	59	123	PHE
32	59	124	GLU
32	59	125	VAL
32	59	127	GLU
32	59	129	THR
32	59	143	GLN
32	59	155	SER
32	59	157	TYR
32	59	164	TYR
32	59	167	GLU
33	69	2	LYS
33	69	4	ILE
33	69	12	LEU
33	69	15	VAL
33	69	33	ARG
33	69	37	VAL
33	69	41	GLU
33	69	56	LYS
33	69	64	GLU
33	69	75	LEU
33	69	76	THR
33	69	77	LEU
33	69	78	THR
33	69	79	ILE
33	69	81	VAL
33	69	82	ARG
33	69	101	LEU
33	69	102	SER
33	69	104	GLN
33	69	105	HIS

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Mol	Chain	Res	Type
33	69	109	ILE
33	69	114	LEU
33	69	117	GLU
33	69	125	GLU
33	69	128	LEU
33	69	131	LYS
33	69	133	HIS
33	69	136	VAL
33	69	142	VAL
34	15	1	MET
34	15	5	VAL
34	15	15	LEU
34	15	16	ILE
34	15	29	LYS
34	15	32	THR
34	15	33	LEU
34	15	34	LEU
34	15	43	THR
34	15	46	VAL
34	15	56	ASN
34	15	63	THR
34	15	65	LYS
34	15	85	ILE
34	15	93	THR
34	15	94	HIS
34	15	99	LEU
34	15	101	HIS
34	15	106	MET
34	15	112	LEU
34	15	127	ASP
34	15	138	LEU
35	25	1	MET
35	25	3	GLN
35	25	12	ASP
35	25	24	VAL
35	25	26	LYS
35	25	32	TYR
35	25	35	VAL
35	25	42	SER
35	25	47	ILE
35	25	49	ARG
35	25	66	LYS

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Mol	Chain	Res	Type
35	25	69	ILE
35	25	70	LYS
35	25	80	ASP
35	25	86	ILE
35	25	94	ARG
35	25	96	THR
35	25	97	ARG
35	25	98	VAL
35	25	113	LYS
35	25	114	ILE
35	25	116	SER
35	25	117	LEU
36	35	6	LEU
36	35	7	ARG
36	35	10	PRO
36	35	14	LYS
36	35	15	ARG
36	35	18	ARG
36	35	21	ARG
36	35	30	THR
36	35	36	LYS
36	35	41	ARG
36	35	50	ARG
36	35	55	ARG
36	35	59	LEU
36	35	61	ARG
36	35	62	LEU
36	35	63	PRO
36	35	65	ARG
36	35	67	MET
36	35	70	GLN
36	35	71	VAL
36	35	75	ILE
36	35	79	ARG
36	35	81	GLN
36	35	85	LEU
36	35	98	GLU
36	35	105	LEU
36	35	110	TYR
36	35	111	ARG
36	35	112	LEU
36	35	114	ILE

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Mol	Chain	Res	Type
36	35	121	LYS
36	35	123	LEU
36	35	124	LYS
36	35	125	VAL
36	35	133	SER
36	35	135	LEU
36	35	138	LEU
36	35	144	GLU
36	35	147	LEU
36	35	149	GLU
37	45	3	MET
37	45	10	ARG
37	45	18	LYS
37	45	25	ASP
37	45	26	TYR
37	45	45	GLN
37	45	51	ARG
37	45	56	ARG
37	45	59	ARG
37	45	60	ARG
37	45	69	PHE
37	45	75	THR
37	45	76	LYS
37	45	79	LEU
37	45	81	VAL
37	45	83	MET
37	45	90	VAL
37	45	91	GLU
37	45	103	MET
37	45	106	VAL
37	45	109	VAL
37	45	110	THR
37	45	116	GLU
37	45	118	LEU
37	45	127	ILE
37	45	134	ARG
37	45	139	GLU
38	55	2	ARG
38	55	18	LEU
38	55	28	LEU
38	55	29	LEU
38	55	34	ILE

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Mol	Chain	Res	Type
38	55	35	THR
38	55	37	THR
38	55	44	LEU
38	55	54	LEU
38	55	56	LYS
38	55	57	ARG
38	55	63	ARG
38	55	65	LEU
38	55	67	LEU
38	55	74	LYS
38	55	75	LEU
38	55	79	LEU
38	55	80	PHE
38	55	81	ASP
38	55	82	GLU
38	55	91	GLN
38	55	97	VAL
39	65	3	ARG
39	65	8	GLU
39	65	12	PHE
39	65	14	VAL
39	65	19	LYS
39	65	21	THR
39	65	27	SER
39	65	30	ARG
39	65	36	TYR
39	65	38	GLN
39	65	52	SER
39	65	54	LEU
39	65	58	LEU
39	65	65	VAL
39	65	82	ILE
39	65	84	GLN
39	65	87	PHE
39	65	93	LYS
39	65	98	VAL
39	65	101	LEU
39	65	106	ARG
39	65	107	GLU
39	65	110	LEU
40	75	1	MET
40	75	6	LEU

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Mol	Chain	Res	Type
40	75	8	LYS
40	75	11	GLU
40	75	13	ARG
40	75	15	VAL
40	75	17	THR
40	75	19	LEU
40	75	27	THR
40	75	28	VAL
40	75	41	ARG
40	75	45	PHE
40	75	55	ASN
40	75	63	VAL
40	75	64	ARG
40	75	67	SER
40	75	74	ARG
40	75	87	ASP
40	75	89	VAL
40	75	91	ARG
40	75	105	LEU
40	75	106	SER
40	75	107	ASP
40	75	112	ARG
40	75	120	ARG
40	75	121	ILE
40	75	124	ASP
40	75	136	GLN
41	85	3	ARG
41	85	5	LYS
41	85	20	LEU
41	85	31	SER
41	85	55	ARG
41	85	57	PHE
41	85	59	ARG
41	85	71	GLN
41	85	74	LEU
41	85	83	LEU
41	85	92	ARG
41	85	97	ASP
41	85	114	LYS
42	95	7	THR
42	95	14	VAL
42	95	18	LEU

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Mol	Chain	Res	Type
42	95	19	LYS
42	95	34	GLU
42	95	35	LEU
42	95	38	LEU
42	95	40	LEU
42	95	44	LYS
42	95	45	THR
42	95	46	VAL
42	95	47	VAL
42	95	49	THR
42	95	57	VAL
42	95	62	LEU
42	95	66	ARG
42	95	71	LEU
42	95	74	LYS
42	95	82	ARG
42	95	83	ARG
42	95	84	LYS
42	95	85	LYS
42	95	89	GLN
42	95	91	TYR
42	95	95	LEU
43	A5	1	MET
43	A5	11	ARG
43	A5	23	LEU
43	A5	37	ARG
43	A5	51	LEU
43	A5	52	GLU
43	A5	65	LEU
43	A5	67	ASP
43	A5	70	TYR
43	A5	92	ARG
43	A5	95	ILE
43	A5	96	ILE
43	A5	100	THR
43	A5	107	LEU
44	B5	3	THR
44	B5	25	LYS
44	B5	27	THR
44	B5	30	VAL
44	B5	36	LYS
44	B5	40	LYS

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Mol	Chain	Res	Type
44	B5	45	THR
44	B5	49	VAL
44	B5	52	VAL
44	B5	63	LYS
44	B5	65	ARG
44	B5	69	TYR
44	B5	70	LEU
44	B5	76	ARG
44	B5	78	LYS
44	B5	81	VAL
44	B5	88	LYS
45	C5	23	ARG
45	C5	24	VAL
45	C5	29	GLU
45	C5	33	LYS
45	C5	38	ILE
45	C5	45	VAL
45	C5	47	LYS
45	C5	50	ARG
45	C5	51	VAL
45	C5	52	SER
45	C5	55	TYR
45	C5	62	GLU
45	C5	63	LYS
45	C5	75	ILE
45	C5	84	ARG
45	C5	85	VAL
45	C5	86	ARG
45	C5	88	LYS
45	C5	89	PHE
45	C5	95	LYS
45	C5	97	ARG
45	C5	98	VAL
45	C5	99	CYS
45	C5	102	CYS
46	D5	11	GLU
46	D5	16	SER
46	D5	19	ARG
46	D5	28	MET
46	D5	53	ILE
46	D5	66	SER
46	D5	70	LEU

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Mol	Chain	Res	Type
46	D5	71	VAL
46	D5	74	VAL
46	D5	76	LEU
46	D5	82	ARG
46	D5	86	VAL
46	D5	87	ASP
46	D5	94	GLU
46	D5	103	ARG
46	D5	107	THR
46	D5	117	LEU
46	D5	121	HIS
46	D5	122	ARG
46	D5	123	ASP
46	D5	133	ILE
46	D5	136	PHE
46	D5	144	LEU
46	D5	157	LEU
46	D5	163	LEU
46	D5	165	VAL
46	D5	170	THR
47	E5	14	ARG
47	E5	29	GLN
47	E5	36	ILE
47	E5	43	THR
47	E5	49	LYS
47	E5	62	LEU
47	E5	64	ASP
47	E5	81	VAL
47	E5	84	LEU
48	F5	20	ARG
48	F5	25	LYS
48	F5	37	ILE
48	F5	38	SER
48	F5	39	LYS
48	F5	41	ARG
48	F5	46	LEU
48	F5	78	LYS
48	F5	82	LEU
48	F5	83	GLU
48	F5	86	SER
48	F5	89	GLU
48	F5	90	ILE

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Mol	Chain	Res	Type
48	F5	91	LYS
48	F5	95	LEU
49	G5	14	ARG
49	G5	16	LEU
49	G5	24	LEU
49	G5	40	SER
49	G5	45	SER
49	G5	47	ASN
49	G5	48	HIS
49	G5	53	LEU
49	G5	59	ARG
50	H5	5	LYS
50	H5	8	LEU
50	H5	18	ASP
50	H5	20	LYS
50	H5	32	GLN
50	H5	38	GLU
50	H5	40	THR
50	H5	44	ARG
50	H5	57	GLU
51	I5	6	HIS
51	I5	9	LEU
51	I5	20	ASN
51	I5	22	ILE
51	I5	25	TYR
51	I5	32	TYR
51	I5	53	GLU
51	I5	59	PHE
51	I5	61	ARG
51	I5	62	ARG
52	J5	3	LYS
52	J5	4	HIS
52	J5	8	LYS
52	J5	15	ARG
52	J5	23	HIS
52	J5	29	THR
52	J5	44	THR
52	J5	48	GLU
52	J5	52	TYR
52	J5	55	ARG
52	J5	56	LYS
53	K5	10	LEU

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Mol	Chain	Res	Type
53	K5	12	GLU
53	K5	23	THR
53	K5	25	LYS
53	K5	27	LYS
53	K5	28	ARG
53	K5	29	ASN
53	K5	36	LEU
53	K5	44	ARG
53	K5	45	LYS
53	K5	47	THR
53	K5	52	VAL
54	L5	2	LYS
54	L5	4	THR
54	L5	8	ASN
54	L5	32	LYS
54	L5	41	ARG
54	L5	43	THR
54	L5	46	VAL
55	M5	4	MET
55	M5	23	VAL
55	M5	25	MET
55	M5	30	ARG
55	M5	34	TRP
55	M5	50	LEU
55	M5	52	LYS
55	M5	53	PRO
55	M5	54	GLU
55	M5	57	ARG
55	M5	59	LYS
55	M5	60	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (29) such sidechains are listed below:

Mol	Chain	Res	Type
3	2E	6	HIS
3	2E	102	ASN
3	2E	136	GLN
6	5E	64	GLN
9	8E	23	ASN
16	7I	14	ASN
19	AI	47	HIS
28	11	96	HIS

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Mol	Chain	Res	Type
28	11	129	ASN
29	21	55	ASN
37	88	13	GLN
38	98	31	HIS
46	H8	34	ASN
2	12	19	HIS
8	72	15	ASN
19	AA	14	HIS
30	39	203	GLN
34	15	56	ASN
36	35	35	HIS
37	45	113	GLN
38	55	13	HIS
40	75	58	ASN
42	95	87	HIS
47	E5	70	GLN
49	G5	48	HIS
50	H5	32	GLN
51	I5	6	HIS
51	I5	20	ASN
53	K5	46	HIS

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	13	1496/1522 (98%)	328 (21%)	30 (2%)
1	1G	1495/1522 (98%)	348 (23%)	39 (2%)
22	1K	74/76 (97%)	31 (41%)	2 (2%)
23	2K	76/77 (98%)	15 (19%)	3 (3%)
23	2L	76/77 (98%)	20 (26%)	3 (3%)
24	1L	75/76 (98%)	26 (34%)	2 (2%)
24	3K	75/76 (98%)	38 (50%)	4 (5%)
24	3L	75/76 (98%)	32 (42%)	2 (2%)
25	4K	12/27 (44%)	4 (33%)	0
25	4L	8/27 (29%)	3 (37%)	1 (12%)
26	14	2908/2917 (99%)	736 (25%)	37 (1%)
26	1H	2911/2917 (99%)	688 (23%)	52 (1%)
27	16	121/122 (99%)	19 (15%)	0
27	1J	121/122 (99%)	31 (25%)	3 (2%)
All	All	9523/9634 (98%)	2319 (24%)	178 (1%)

All (2319) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	13	5	U
1	13	6	G
1	13	7	G
1	13	8	A
1	13	9	G
1	13	21	G
1	13	32	A
1	13	39	G
1	13	47	C
1	13	48	C
1	13	50	A
1	13	51	A
1	13	54	C
1	13	61	G
1	13	65	U
1	13	66	G
1	13	77	C
1	13	78	G
1	13	91	C
1	13	101	A
1	13	108	G
1	13	116	A
1	13	119	A
1	13	121	C
1	13	129(A)	G
1	13	130	A
1	13	131	C
1	13	132	C
1	13	137	C
1	13	142	G
1	13	144	G
1	13	150	C
1	13	151	A
1	13	160	A
1	13	161	A
1	13	163	C
1	13	169	C
1	13	172	A
1	13	173	U
1	13	174	C
1	13	186(A)	C
1	13	188	U
1	13	189	U

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Mol	Chain	Res	Type
1	13	190	G
1	13	191(A)	G
1	13	195	A
1	13	197	A
1	13	199	G
1	13	201	C
1	13	208	U
1	13	209	U
1	13	210	U
1	13	216	G
1	13	222	U
1	13	226	G
1	13	231	G
1	13	233	C
1	13	243	A
1	13	244	U
1	13	245	C
1	13	247	G
1	13	251	G
1	13	262	A
1	13	266	G
1	13	267	C
1	13	273	A
1	13	281	G
1	13	289	G
1	13	302	G
1	13	306	G
1	13	311	C
1	13	321	A
1	13	324	G
1	13	328	C
1	13	329	A
1	13	332	G
1	13	339	C
1	13	344	A
1	13	345	C
1	13	346	G
1	13	347	G
1	13	352	C
1	13	353	A
1	13	354	G
1	13	357	G

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Mol	Chain	Res	Type
1	13	365	U
1	13	367	U
1	13	372	C
1	13	384	G
1	13	388	G
1	13	390	C
1	13	397	A
1	13	398	C
1	13	406	G
1	13	412	A
1	13	413	G
1	13	414	A
1	13	419	C
1	13	421	U
1	13	423	G
1	13	424	G
1	13	428	G
1	13	429	U
1	13	455	C
1	13	465	A
1	13	466	C
1	13	467	G
1	13	485	G
1	13	496	A
1	13	497	U
1	13	505	G
1	13	509	A
1	13	510	A
1	13	511	C
1	13	513	C
1	13	518	C
1	13	524	G
1	13	527	G
1	13	531	U
1	13	532	A
1	13	533	A
1	13	536	C
1	13	545	C
1	13	547	A
1	13	549	C
1	13	559	A
1	13	560	U

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Mol	Chain	Res	Type
1	13	561	U
1	13	571	U
1	13	572	A
1	13	573	A
1	13	576	G
1	13	577	G
1	13	607	A
1	13	610	G
1	13	620	C
1	13	630	G
1	13	631	G
1	13	632	A
1	13	633	G
1	13	639	G
1	13	650	G
1	13	653	A
1	13	655	A
1	13	656	C
1	13	665	A
1	13	666	G
1	13	680	C
1	13	687	A
1	13	688	G
1	13	702	A
1	13	704	A
1	13	723	U
1	13	724	G
1	13	734	G
1	13	748	C
1	13	749	C
1	13	750	G
1	13	753	A
1	13	755	G
1	13	777	A
1	13	787	A
1	13	792	A
1	13	793	U
1	13	794	A
1	13	805	C
1	13	813	U
1	13	815	A
1	13	817	C

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Mol	Chain	Res	Type
1	13	818	G
1	13	827	U
1	13	828	A
1	13	836	G
1	13	841	U
1	13	842	C
1	13	843	U
1	13	848	C
1	13	853	G
1	13	859	A
1	13	864	A
1	13	870	U
1	13	874	G
1	13	888	G
1	13	902	G
1	13	905	U
1	13	914	A
1	13	916	G
1	13	922	G
1	13	926	G
1	13	927	G
1	13	934	C
1	13	935	A
1	13	936	C
1	13	940	C
1	13	955	U
1	13	960	U
1	13	968	A
1	13	969	A
1	13	971	G
1	13	974	A
1	13	975	A
1	13	976	G
1	13	977	A
1	13	982	U
1	13	983	A
1	13	991	U
1	13	992	U
1	13	993	G
1	13	1004	A
1	13	1006	C
1	13	1007	C

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Mol	Chain	Res	Type
1	13	1008	C
1	13	1009	G
1	13	1012	U
1	13	1017	G
1	13	1021	G
1	13	1024	G
1	13	1025	U
1	13	1028	C
1	13	1029	G
1	13	1030	C
1	13	1032(A)	G
1	13	1033	G
1	13	1040	U
1	13	1046	A
1	13	1049	U
1	13	1053	G
1	13	1054	C
1	13	1064	G
1	13	1065	U
1	13	1066	C
1	13	1081	G
1	13	1082	G
1	13	1094	G
1	13	1095	U
1	13	1101	A
1	13	1120	G
1	13	1122	U
1	13	1124	G
1	13	1125	U
1	13	1126	U
1	13	1127	G
1	13	1129	C
1	13	1130	A
1	13	1131	G
1	13	1136	U
1	13	1137	C
1	13	1138	G
1	13	1139	G
1	13	1146	A
1	13	1147	C
1	13	1152	A
1	13	1154	G

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Mol	Chain	Res	Type
1	13	1157	A
1	13	1158	C
1	13	1159	U
1	13	1161	C
1	13	1177	G
1	13	1178	G
1	13	1181	G
1	13	1191	A
1	13	1196	U
1	13	1197	G
1	13	1201	A
1	13	1212	U
1	13	1213	A
1	13	1225	A
1	13	1227	A
1	13	1238	A
1	13	1240	U
1	13	1241	G
1	13	1250	A
1	13	1253	G
1	13	1256	A
1	13	1257	U
1	13	1258	G
1	13	1270	C
1	13	1273	G
1	13	1275	A
1	13	1278	U
1	13	1280	A
1	13	1281	U
1	13	1282	C
1	13	1286	A
1	13	1287	A
1	13	1299	A
1	13	1300	G
1	13	1302	U
1	13	1320	C
1	13	1322	C
1	13	1323	G
1	13	1331	G
1	13	1335	C
1	13	1336	C
1	13	1337	G

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Mol	Chain	Res	Type
1	13	1338	G
1	13	1340	A
1	13	1346	A
1	13	1347	G
1	13	1350	A
1	13	1353	G
1	13	1368	G
1	13	1370	G
1	13	1377	A
1	13	1398	A
1	13	1401	G
1	13	1419	G
1	13	1422	G
1	13	1442	G
1	13	1443	G
1	13	1446	A
1	13	1451	A
1	13	1452	C
1	13	1453	G
1	13	1467	G
1	13	1469	G
1	13	1487	G
1	13	1492	A
1	13	1497	G
1	13	1499	A
1	13	1502	A
1	13	1503	A
1	13	1504	G
1	13	1505	G
1	13	1506	U
1	13	1517	G
1	13	1529	G
1	13	1530	G
22	1K	2	C
22	1K	9	A
22	1K	10	G
22	1K	11	C
22	1K	14	A
22	1K	16	U
22	1K	17	C
22	1K	18	G
22	1K	19	G

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Mol	Chain	Res	Type
22	1K	21	A
22	1K	22	G
22	1K	24	G
22	1K	25	C
22	1K	26	A
22	1K	28	G
22	1K	41	C
22	1K	42	C
22	1K	44	G
22	1K	45	U
22	1K	47	U
22	1K	49	C
22	1K	52	G
22	1K	55	PSU
22	1K	61	C
22	1K	64	A
22	1K	68	C
22	1K	70	G
22	1K	73	A
22	1K	74	C
22	1K	75	C
22	1K	76	A
23	2K	2	G
23	2K	8	4SU
23	2K	9	G
23	2K	18	C
23	2K	20	G
23	2K	21	U
23	2K	22	A
23	2K	23	G
23	2K	31	G
23	2K	44	A
23	2K	48	U
23	2K	49	C
23	2K	50	G
23	2K	57	C
23	2K	77	A
24	3K	2	C
24	3K	3	C
24	3K	7	A
24	3K	9	A
24	3K	10	G

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Mol	Chain	Res	Type
24	3K	13	C
24	3K	14	A
24	3K	17	C
24	3K	19	G
24	3K	20	U
24	3K	21	A
24	3K	22	G
24	3K	31	A
24	3K	35	A
24	3K	36	A
24	3K	39	U
24	3K	40	C
24	3K	41	C
24	3K	42	C
24	3K	45	U
24	3K	46	G
24	3K	47	U
24	3K	48	C
24	3K	49	C
24	3K	52	G
24	3K	55	U
24	3K	56	C
24	3K	58	A
24	3K	59	U
24	3K	63	G
24	3K	64	A
24	3K	65	G
24	3K	66	U
24	3K	67	C
24	3K	70	G
24	3K	72	C
24	3K	73	A
24	3K	76	A
25	4K	14	A
25	4K	15	A
25	4K	21	C
25	4K	25	A
26	1H	5	A
26	1H	9	U
26	1H	12	U
26	1H	15	G
26	1H	17	G

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Mol	Chain	Res	Type
26	1H	27	G
26	1H	34	C
26	1H	35	G
26	1H	36	G
26	1H	39	C
26	1H	46	C
26	1H	51	G
26	1H	63	U
26	1H	64	A
26	1H	71	A
26	1H	72	U
26	1H	74	A
26	1H	75	G
26	1H	85	G
26	1H	95	G
26	1H	102	G
26	1H	112	U
26	1H	118	A
26	1H	119	A
26	1H	120	U
26	1H	124	G
26	1H	125	G
26	1H	155	C
26	1H	163	U
26	1H	164	U
26	1H	171	G
26	1H	175	G
26	1H	181	A
26	1H	188	G
26	1H	196	A
26	1H	199	A
26	1H	201	C
26	1H	215	G
26	1H	216	A
26	1H	222	A
26	1H	223	A
26	1H	224	G
26	1H	228	A
26	1H	229	A
26	1H	230	U
26	1H	233	A
26	1H	244	A

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Mol	Chain	Res	Type
26	1H	245	G
26	1H	248	G
26	1H	250	G
26	1H	252	G
26	1H	261	G
26	1H	264	C
26	1H	269	U
26	1H	270(D)	C
26	1H	270(F)	U
26	1H	270(K)	C
26	1H	270(M)	U
26	1H	270(N)	G
26	1H	270(P)	C
26	1H	271(C)	U
26	1H	271	G
26	1H	273(E)	U
26	1H	274	G
26	1H	275	G
26	1H	277	C
26	1H	278	A
26	1H	299	A
26	1H	308	G
26	1H	311	A
26	1H	315	G
26	1H	323	G
26	1H	324	A
26	1H	327	G
26	1H	329	G
26	1H	330	A
26	1H	331	A
26	1H	342	G
26	1H	352	G
26	1H	363	G
26	1H	363(B)	G
26	1H	364	C
26	1H	372	G
26	1H	382	G
26	1H	386	G
26	1H	389	G
26	1H	396	G
26	1H	405	U
26	1H	411	G

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Mol	Chain	Res	Type
26	1H	412	A
26	1H	427	U
26	1H	428	A
26	1H	443	A
26	1H	444	C
26	1H	447	A
26	1H	448	U
26	1H	452	G
26	1H	454	A
26	1H	455	C
26	1H	457	A
26	1H	459	U
26	1H	460	A
26	1H	470	A
26	1H	471	A
26	1H	481	G
26	1H	482	A
26	1H	488	G
26	1H	501	A
26	1H	502	A
26	1H	505	A
26	1H	508	G
26	1H	509	C
26	1H	529	A
26	1H	531	C
26	1H	532	A
26	1H	533	G
26	1H	545	G
26	1H	546	C
26	1H	549	G
26	1H	556	G
26	1H	559	G
26	1H	563	G
26	1H	564	C
26	1H	567	A
26	1H	573	G
26	1H	575	A
26	1H	577	G
26	1H	586	A
26	1H	588	U
26	1H	603	A
26	1H	607	U

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Mol	Chain	Res	Type
26	1H	614	U
26	1H	615	G
26	1H	617	G
26	1H	621	A
26	1H	622	G
26	1H	627	A
26	1H	632	A
26	1H	636	G
26	1H	637	A
26	1H	645	C
26	1H	646	A
26	1H	647	G
26	1H	654	A
26	1H	654(A)	A
26	1H	654(G)	C
26	1H	654(I)	C
26	1H	654(K)	C
26	1H	654(M)	C
26	1H	654(O)	G
26	1H	654(T)	A
26	1H	654(V)	A
26	1H	664	C
26	1H	669	G
26	1H	672	C
26	1H	686	G
26	1H	699	A
26	1H	717	G
26	1H	724	U
26	1H	729	G
26	1H	730	C
26	1H	731	C
26	1H	745	G
26	1H	752	A
26	1H	753	C
26	1H	764	A
26	1H	765	G
26	1H	775	G
26	1H	776	G
26	1H	777	A
26	1H	782	A
26	1H	784	A
26	1H	785	G

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Mol	Chain	Res	Type
26	1H	790	C
26	1H	792	G
26	1H	801	G
26	1H	805	G
26	1H	812	C
26	1H	824	A
26	1H	827	U
26	1H	828	U
26	1H	832	G
26	1H	845	G
26	1H	846	C
26	1H	847	U
26	1H	859	G
26	1H	861	A
26	1H	866	A
26	1H	870	A
26	1H	879	G
26	1H	880	G
26	1H	881	G
26	1H	882	G
26	1H	884	C
26	1H	885	C
26	1H	886	C
26	1H	887	A
26	1H	888	C
26	1H	890	A
26	1H	892	G
26	1H	893	C
26	1H	894	C
26	1H	895	U
26	1H	896	A
26	1H	897	C
26	1H	900	A
26	1H	901	A
26	1H	910	A
26	1H	914	C
26	1H	917	A
26	1H	918	A
26	1H	932	G
26	1H	938	G
26	1H	940	G
26	1H	941	A

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Mol	Chain	Res	Type
26	1H	946	G
26	1H	947	G
26	1H	952	G
26	1H	953	A
26	1H	959	A
26	1H	961	C
26	1H	974	G
26	1H	974(A)	C
26	1H	975	G
26	1H	981	A
26	1H	982	C
26	1H	983	A
26	1H	996	A
26	1H	997	G
26	1H	1003	G
26	1H	1005	C
26	1H	1011	G
26	1H	1012	U
26	1H	1013	C
26	1H	1016	G
26	1H	1020	A
26	1H	1022	G
26	1H	1023	U
26	1H	1025	G
26	1H	1026	U
26	1H	1027	A
26	1H	1033	U
26	1H	1046	A
26	1H	1047	G
26	1H	1054	A
26	1H	1057	A
26	1H	1060	U
26	1H	1061	U
26	1H	1062	G
26	1H	1064	C
26	1H	1068	G
26	1H	1070	A
26	1H	1071	G
26	1H	1072	C
26	1H	1073	A
26	1H	1076	C
26	1H	1078	U

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Mol	Chain	Res	Type
26	1H	1079	C
26	1H	1082	U
26	1H	1084	A
26	1H	1085	A
26	1H	1086	A
26	1H	1087	G
26	1H	1088	A
26	1H	1090	U
26	1H	1095	A
26	1H	1096	A
26	1H	1097	U
26	1H	1104	C
26	1H	1106	G
26	1H	1110	G
26	1H	1111	A
26	1H	1112	G
26	1H	1121	C
26	1H	1122	G
26	1H	1126	A
26	1H	1127	A
26	1H	1128	A
26	1H	1129	A
26	1H	1130	U
26	1H	1135	C
26	1H	1136	G
26	1H	1139	G
26	1H	1142	U
26	1H	1142(A)	A
26	1H	1144	G
26	1H	1155	A
26	1H	1156	A
26	1H	1171	G
26	1H	1174	A
26	1H	1175	U
26	1H	1176	G
26	1H	1178	C
26	1H	1179	C
26	1H	1180	C
26	1H	1190	G
26	1H	1194	A
26	1H	1195	G
26	1H	1204	A

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Mol	Chain	Res	Type
26	1H	1205	U
26	1H	1220	A
26	1H	1221	C
26	1H	1229(A)	G
26	1H	1244	G
26	1H	1253	A
26	1H	1255	U
26	1H	1256	G
26	1H	1265	A
26	1H	1267	U
26	1H	1271	G
26	1H	1272	A
26	1H	1273	U
26	1H	1274	A
26	1H	1280	G
26	1H	1292	U
26	1H	1300	U
26	1H	1301	A
26	1H	1305	C
26	1H	1313	U
26	1H	1314	C
26	1H	1321	A
26	1H	1329	U
26	1H	1332	G
26	1H	1344	G
26	1H	1345	C
26	1H	1349	A
26	1H	1352	U
26	1H	1358	G
26	1H	1359	A
26	1H	1360	A
26	1H	1365	A
26	1H	1368	G
26	1H	1369	G
26	1H	1379	A
26	1H	1380	G
26	1H	1384	A
26	1H	1385	G
26	1H	1386	C
26	1H	1388	G
26	1H	1389	G
26	1H	1390	U

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Mol	Chain	Res	Type
26	1H	1416	G
26	1H	1417	C
26	1H	1420	U
26	1H	1421	G
26	1H	1427	A
26	1H	1428	C
26	1H	1431	U
26	1H	1437	C
26	1H	1444(A)	A
26	1H	1449	A
26	1H	1449(A)	G
26	1H	1455	G
26	1H	1458	C
26	1H	1459	G
26	1H	1460	A
26	1H	1461	G
26	1H	1467	C
26	1H	1471	A
26	1H	1473	G
26	1H	1483	G
26	1H	1492	G
26	1H	1493	C
26	1H	1494	A
26	1H	1497	U
26	1H	1506	C
26	1H	1507	A
26	1H	1508	A
26	1H	1509	C
26	1H	1510	A
26	1H	1511	A
26	1H	1517	G
26	1H	1520	U
26	1H	1526	G
26	1H	1534	G
26	1H	1535	U
26	1H	1536	A
26	1H	1537	C
26	1H	1538	G
26	1H	1540	G
26	1H	1543	A
26	1H	1544	C
26	1H	1545	A

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Mol	Chain	Res	Type
26	1H	1548	C
26	1H	1554	A
26	1H	1558	A
26	1H	1559	G
26	1H	1560	G
26	1H	1562	A
26	1H	1566	A
26	1H	1569	A
26	1H	1578	U
26	1H	1580	A
26	1H	1586	A
26	1H	1592	C
26	1H	1608	A
26	1H	1609	A
26	1H	1610	A
26	1H	1617	C
26	1H	1618	A
26	1H	1634	A
26	1H	1647	G
26	1H	1648	C
26	1H	1651	G
26	1H	1654	A
26	1H	1669	A
26	1H	1674	G
26	1H	1684	C
26	1H	1694	C
26	1H	1695	G
26	1H	1699	G
26	1H	1728	G
26	1H	1729	A
26	1H	1730	U
26	1H	1731	G
26	1H	1756	G
26	1H	1758	G
26	1H	1762	A
26	1H	1763	G
26	1H	1764	G
26	1H	1772	G
26	1H	1773	A
26	1H	1782	C
26	1H	1787	A
26	1H	1791	A

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Mol	Chain	Res	Type
26	1H	1799	G
26	1H	1800	C
26	1H	1801	G
26	1H	1802	A
26	1H	1811	G
26	1H	1816	G
26	1H	1817	G
26	1H	1829	A
26	1H	1835	G
26	1H	1839	G
26	1H	1847	A
26	1H	1858	G
26	1H	1870	C
26	1H	1878	G
26	1H	1889	A
26	1H	1896	G
26	1H	1900	A
26	1H	1906	G
26	1H	1914	C
26	1H	1919	A
26	1H	1920	C
26	1H	1926	U
26	1H	1929	G
26	1H	1930	G
26	1H	1931	U
26	1H	1938	A
26	1H	1940	U
26	1H	1941	C
26	1H	1951	U
26	1H	1952	A
26	1H	1955	U
26	1H	1960	A
26	1H	1967	C
26	1H	1969	A
26	1H	1970	A
26	1H	1971	A
26	1H	1972	A
26	1H	1982	C
26	1H	1993	U
26	1H	2005	A
26	1H	2020	A
26	1H	2021	C

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Mol	Chain	Res	Type
26	1H	2023	G
26	1H	2030	A
26	1H	2031	A
26	1H	2033	A
26	1H	2036	C
26	1H	2043	C
26	1H	2046	G
26	1H	2051	A
26	1H	2055	C
26	1H	2056	G
26	1H	2059	A
26	1H	2060	A
26	1H	2061	G
26	1H	2062	A
26	1H	2063	C
26	1H	2069	G
26	1H	2078	C
26	1H	2080	G
26	1H	2096	U
26	1H	2099	U
26	1H	2108	C
26	1H	2111	C
26	1H	2112	G
26	1H	2113	U
26	1H	2114	A
26	1H	2116	G
26	1H	2118	U
26	1H	2122	U
26	1H	2126	A
26	1H	2127	G
26	1H	2128	C
26	1H	2131	G
26	1H	2132	U
26	1H	2133	G
26	1H	2135	A
26	1H	2136	C
26	1H	2139	C
26	1H	2147	G
26	1H	2148	G
26	1H	2154	G
26	1H	2157	G
26	1H	2158	A

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Mol	Chain	Res	Type
26	1H	2161	C
26	1H	2166	G
26	1H	2167	U
26	1H	2168	G
26	1H	2170	A
26	1H	2171	A
26	1H	2172	U
26	1H	2173	A
26	1H	2174	C
26	1H	2176	A
26	1H	2181	G
26	1H	2190	G
26	1H	2192	G
26	1H	2198	A
26	1H	2205	C
26	1H	2210	G
26	1H	2212	A
26	1H	2213	U
26	1H	2215	G
26	1H	2225	A
26	1H	2226	C
26	1H	2236	C
26	1H	2238	G
26	1H	2240	C
26	1H	2268	A
26	1H	2269	A
26	1H	2273	A
26	1H	2275	C
26	1H	2278	A
26	1H	2280	G
26	1H	2281	C
26	1H	2283	C
26	1H	2287	A
26	1H	2288	A
26	1H	2297	C
26	1H	2305	A
26	1H	2307	G
26	1H	2308	G
26	1H	2310	A
26	1H	2311	A
26	1H	2312	U
26	1H	2315	G

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Mol	Chain	Res	Type
26	1H	2319	G
26	1H	2320	A
26	1H	2324	C
26	1H	2325	G
26	1H	2326	C
26	1H	2327	A
26	1H	2334	G
26	1H	2335	A
26	1H	2336	A
26	1H	2342	C
26	1H	2346	A
26	1H	2347	C
26	1H	2350	C
26	1H	2360	A
26	1H	2372	G
26	1H	2376	A
26	1H	2377	A
26	1H	2379	G
26	1H	2383	G
26	1H	2385	C
26	1H	2392	A
26	1H	2393	A
26	1H	2395	C
26	1H	2398	U
26	1H	2402	C
26	1H	2403	C
26	1H	2405	G
26	1H	2406	U
26	1H	2410	G
26	1H	2414	G
26	1H	2418	A
26	1H	2424	C
26	1H	2425	A
26	1H	2428	G
26	1H	2429	G
26	1H	2430	A
26	1H	2439	A
26	1H	2440	C
26	1H	2441	C
26	1H	2448	A
26	1H	2450	A
26	1H	2469	A

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Mol	Chain	Res	Type
26	1H	2470	G
26	1H	2476	A
26	1H	2478	A
26	1H	2482	G
26	1H	2484	G
26	1H	2487	G
26	1H	2497	A
26	1H	2502	G
26	1H	2505	G
26	1H	2506	U
26	1H	2518	A
26	1H	2520	C
26	1H	2529	G
26	1H	2549	G
26	1H	2554	U
26	1H	2566	A
26	1H	2567	G
26	1H	2573	C
26	1H	2576	G
26	1H	2582	G
26	1H	2585	U
26	1H	2601	C
26	1H	2602	A
26	1H	2609	U
26	1H	2610	C
26	1H	2611	U
26	1H	2612	C
26	1H	2615	U
26	1H	2629	A
26	1H	2636	U
26	1H	2641	G
26	1H	2654	A
26	1H	2660	A
26	1H	2665	A
26	1H	2666	C
26	1H	2673	G
26	1H	2689	U
26	1H	2690	C
26	1H	2691	C
26	1H	2698	U
26	1H	2700	C
26	1H	2701	C

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Mol	Chain	Res	Type
26	1H	2702	U
26	1H	2703	C
26	1H	2704	C
26	1H	2705	A
26	1H	2707	G
26	1H	2712(A)	A
26	1H	2713	A
26	1H	2714	G
26	1H	2718	G
26	1H	2719	G
26	1H	2721	A
26	1H	2726	U
26	1H	2733	A
26	1H	2756	U
26	1H	2757	A
26	1H	2758	A
26	1H	2764	A
26	1H	2765	A
26	1H	2766	G
26	1H	2778	A
26	1H	2779	U
26	1H	2780	G
26	1H	2781	A
26	1H	2789	C
26	1H	2790	A
26	1H	2791	C
26	1H	2793	G
26	1H	2794	C
26	1H	2795	G
26	1H	2797	U
26	1H	2798	C
26	1H	2801	A
26	1H	2802	G
26	1H	2808	U
26	1H	2818	G
26	1H	2820	A
26	1H	2821	A
26	1H	2832	U
26	1H	2833	G
26	1H	2834	G
26	1H	2835	A
26	1H	2864	G

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Mol	Chain	Res	Type
26	1H	2871	C
26	1H	2872	G
26	1H	2876	G
26	1H	2883	A
26	1H	2891	G
26	1H	2892	A
26	1H	2893	G
26	1H	2894	G
26	1H	2895	U
26	1H	2899	G
26	1H	2902	C
27	16	7	G
27	16	9	G
27	16	13	A
27	16	15	A
27	16	16	G
27	16	25	A
27	16	33	G
27	16	39	A
27	16	40	U
27	16	41	U
27	16	45	A
27	16	56	G
27	16	66	A
27	16	73	A
27	16	81	G
27	16	82	G
27	16	105	G
27	16	109	G
27	16	115	G
1	1G	5	U
1	1G	7	G
1	1G	9	G
1	1G	16	A
1	1G	22	G
1	1G	26	A
1	1G	31	G
1	1G	32	A
1	1G	39	G
1	1G	41	G
1	1G	47	C
1	1G	48	C

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Mol	Chain	Res	Type
1	1G	50	A
1	1G	51	A
1	1G	54	C
1	1G	65	U
1	1G	76	G
1	1G	79	G
1	1G	90	C
1	1G	91	C
1	1G	101	A
1	1G	115	G
1	1G	116	A
1	1G	121	C
1	1G	129(A)	G
1	1G	131	C
1	1G	132	C
1	1G	144	G
1	1G	146	G
1	1G	154	C
1	1G	163	C
1	1G	167	G
1	1G	173	U
1	1G	174	C
1	1G	182	U
1	1G	185	A
1	1G	187	C
1	1G	188	U
1	1G	189	U
1	1G	190	G
1	1G	191(A)	G
1	1G	191(D)	U
1	1G	195	A
1	1G	197	A
1	1G	198	G
1	1G	209	U
1	1G	210	U
1	1G	216	G
1	1G	238	G
1	1G	242	C
1	1G	244	U
1	1G	247	G
1	1G	250	A
1	1G	251	G

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Mol	Chain	Res	Type
1	1G	266	G
1	1G	267	C
1	1G	274	A
1	1G	281	G
1	1G	283	C
1	1G	289	G
1	1G	290	C
1	1G	298	A
1	1G	321	A
1	1G	328	C
1	1G	329	A
1	1G	332	G
1	1G	346	G
1	1G	349	A
1	1G	350	G
1	1G	351	G
1	1G	352	C
1	1G	353	A
1	1G	354	G
1	1G	356	A
1	1G	366	C
1	1G	367	U
1	1G	372	C
1	1G	388	G
1	1G	397	A
1	1G	398	C
1	1G	406	G
1	1G	411	A
1	1G	412	A
1	1G	413	G
1	1G	414	A
1	1G	421	U
1	1G	422	C
1	1G	423	G
1	1G	424	G
1	1G	429	U
1	1G	430	A
1	1G	439	A
1	1G	442	C
1	1G	448	A
1	1G	452	A
1	1G	465	A

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Mol	Chain	Res	Type
1	1G	466	C
1	1G	467	G
1	1G	475	G
1	1G	478	A
1	1G	484	G
1	1G	485	G
1	1G	486	U
1	1G	496	A
1	1G	497	U
1	1G	502	G
1	1G	505	G
1	1G	509	A
1	1G	510	A
1	1G	511	C
1	1G	512	U
1	1G	513	C
1	1G	518	C
1	1G	521	G
1	1G	524	G
1	1G	527	G
1	1G	529	G
1	1G	530	G
1	1G	531	U
1	1G	532	A
1	1G	533	A
1	1G	534	U
1	1G	536	C
1	1G	546	G
1	1G	547	A
1	1G	559	A
1	1G	561	U
1	1G	564	C
1	1G	566	G
1	1G	572	A
1	1G	573	A
1	1G	575	G
1	1G	576	G
1	1G	577	G
1	1G	581	G
1	1G	607	A
1	1G	611	A
1	1G	614	A

Continued on next page...

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Mol	Chain	Res	Type
1	1G	618	C
1	1G	629	G
1	1G	630	G
1	1G	632	A
1	1G	633	G
1	1G	652	U
1	1G	653	A
1	1G	661	G
1	1G	665	A
1	1G	687	A
1	1G	688	G
1	1G	702	A
1	1G	703	G
1	1G	721	G
1	1G	724	G
1	1G	731	G
1	1G	749	C
1	1G	760	G
1	1G	769	G
1	1G	777	A
1	1G	778	G
1	1G	782	A
1	1G	792	A
1	1G	794	A
1	1G	803	G
1	1G	817	C
1	1G	821	G
1	1G	827	U
1	1G	828	A
1	1G	842	C
1	1G	843	U
1	1G	848	C
1	1G	851	G
1	1G	859	A
1	1G	867	G
1	1G	873	A
1	1G	885	G
1	1G	887	G
1	1G	914	A
1	1G	916	G
1	1G	926	G
1	1G	927	G

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Mol	Chain	Res	Type
1	1G	934	C
1	1G	935	A
1	1G	936	C
1	1G	958	A
1	1G	960	U
1	1G	961	U
1	1G	966	G
1	1G	968	A
1	1G	969	A
1	1G	971	G
1	1G	972	C
1	1G	974	A
1	1G	975	A
1	1G	976	G
1	1G	977	A
1	1G	978	A
1	1G	980	C
1	1G	981	U
1	1G	983	A
1	1G	991	U
1	1G	992	U
1	1G	993	G
1	1G	1004	A
1	1G	1006	C
1	1G	1009	G
1	1G	1024	G
1	1G	1025	U
1	1G	1027	C
1	1G	1028	C
1	1G	1028(B)	C
1	1G	1029	G
1	1G	1031	G
1	1G	1032(A)	G
1	1G	1033	G
1	1G	1035	A
1	1G	1036	G
1	1G	1040	U
1	1G	1046	A
1	1G	1047	G
1	1G	1050	G
1	1G	1051	C
1	1G	1053	G

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Mol	Chain	Res	Type
1	1G	1054	C
1	1G	1055	A
1	1G	1081	G
1	1G	1082	G
1	1G	1085	U
1	1G	1086	U
1	1G	1094	G
1	1G	1095	U
1	1G	1101	A
1	1G	1118	C
1	1G	1123	A
1	1G	1125	U
1	1G	1127	G
1	1G	1128	C
1	1G	1129	C
1	1G	1131	G
1	1G	1135	U
1	1G	1137	C
1	1G	1138	G
1	1G	1139	G
1	1G	1145	C
1	1G	1146	A
1	1G	1147	C
1	1G	1154	G
1	1G	1157	A
1	1G	1158	C
1	1G	1159	U
1	1G	1160	G
1	1G	1171	G
1	1G	1177	G
1	1G	1178	G
1	1G	1181	G
1	1G	1182	G
1	1G	1183	A
1	1G	1186	G
1	1G	1187	G
1	1G	1188	A
1	1G	1190	G
1	1G	1193	G
1	1G	1196	U
1	1G	1198	G
1	1G	1200	C

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Mol	Chain	Res	Type
1	1G	1201	A
1	1G	1202	G
1	1G	1211	U
1	1G	1212	U
1	1G	1213	A
1	1G	1214	C
1	1G	1225	A
1	1G	1232	U
1	1G	1233	G
1	1G	1238	A
1	1G	1240	U
1	1G	1241	G
1	1G	1256	A
1	1G	1257	U
1	1G	1258	G
1	1G	1260	C
1	1G	1263	C
1	1G	1267	C
1	1G	1269	A
1	1G	1274	G
1	1G	1275	A
1	1G	1278	U
1	1G	1279	A
1	1G	1280	A
1	1G	1281	U
1	1G	1286	A
1	1G	1287	A
1	1G	1288	A
1	1G	1290	G
1	1G	1293	G
1	1G	1297	C
1	1G	1298	C
1	1G	1299	A
1	1G	1301	U
1	1G	1303	C
1	1G	1305	G
1	1G	1313	U
1	1G	1322	C
1	1G	1323	G
1	1G	1331	G
1	1G	1335	C
1	1G	1336	C

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Mol	Chain	Res	Type
1	1G	1346	A
1	1G	1347	G
1	1G	1353	G
1	1G	1362(A)	C
1	1G	1363	A
1	1G	1364	U
1	1G	1368	G
1	1G	1370	G
1	1G	1379	G
1	1G	1392	G
1	1G	1396	A
1	1G	1397	C
1	1G	1398	A
1	1G	1401	G
1	1G	1402	C
1	1G	1406	U
1	1G	1419	G
1	1G	1442	G
1	1G	1443	G
1	1G	1446	A
1	1G	1449	C
1	1G	1450	U
1	1G	1451	A
1	1G	1452	C
1	1G	1453	G
1	1G	1454	G
1	1G	1469	G
1	1G	1482	G
1	1G	1491	G
1	1G	1492	A
1	1G	1494	G
1	1G	1499	A
1	1G	1502	A
1	1G	1503	A
1	1G	1504	G
1	1G	1506	U
1	1G	1517	G
1	1G	1519	A
1	1G	1520	G
1	1G	1525	G
1	1G	1529	G
1	1G	1530	G

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Mol	Chain	Res	Type
24	1L	2	C
24	1L	8	U
24	1L	9	A
24	1L	10	G
24	1L	11	C
24	1L	16	U
24	1L	17	C
24	1L	19	G
24	1L	21	A
24	1L	22	G
24	1L	25	C
24	1L	27	G
24	1L	36	A
24	1L	40	C
24	1L	41	C
24	1L	46	G
24	1L	47	U
24	1L	48	C
24	1L	59	U
24	1L	61	C
24	1L	69	G
24	1L	70	G
24	1L	73	A
24	1L	74	C
24	1L	75	C
24	1L	76	A
23	2L	2	G
23	2L	8	4SU
23	2L	9	G
23	2L	13	C
23	2L	16	C
23	2L	17	C
23	2L	19	G
23	2L	20	G
23	2L	21	U
23	2L	22	A
23	2L	23	G
23	2L	27	G
23	2L	32	G
23	2L	48	U
23	2L	49	C
23	2L	50	G

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Mol	Chain	Res	Type
23	2L	53	G
23	2L	57	C
23	2L	61	U
23	2L	68	C
24	3L	2	C
24	3L	9	A
24	3L	13	C
24	3L	16	U
24	3L	17	C
24	3L	18	G
24	3L	19	G
24	3L	20	U
24	3L	21	A
24	3L	22	G
24	3L	23	A
24	3L	26	A
24	3L	31	A
24	3L	33	U
24	3L	36	A
24	3L	37	A
24	3L	38	A
24	3L	40	C
24	3L	44	G
24	3L	46	G
24	3L	47	U
24	3L	48	C
24	3L	54	U
24	3L	58	A
24	3L	59	U
24	3L	61	C
24	3L	62	C
24	3L	65	G
24	3L	66	U
24	3L	72	C
24	3L	73	A
24	3L	76	A
25	4L	13	A
25	4L	14	A
25	4L	19	U
26	14	2	G
26	14	3	U
26	14	4	C

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Mol	Chain	Res	Type
26	14	5	A
26	14	6	A
26	14	9	U
26	14	14	A
26	14	15	G
26	14	34	C
26	14	35	G
26	14	46	C
26	14	54	G
26	14	55	G
26	14	58	G
26	14	71	A
26	14	72	U
26	14	74	A
26	14	75	G
26	14	82	G
26	14	84	A
26	14	91	A
26	14	93	C
26	14	95	G
26	14	99	U
26	14	102	G
26	14	118	A
26	14	119	A
26	14	120	U
26	14	121	G
26	14	125	G
26	14	129	C
26	14	131	G
26	14	138	G
26	14	139	G
26	14	140	A
26	14	154	G
26	14	155	C
26	14	161	U
26	14	162	U
26	14	173	G
26	14	174	C
26	14	181	A
26	14	182	A
26	14	196	A
26	14	199	A

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Mol	Chain	Res	Type
26	14	201	C
26	14	205	G
26	14	214	G
26	14	215	G
26	14	216	A
26	14	217	G
26	14	221	A
26	14	222	A
26	14	225	A
26	14	229	A
26	14	232	G
26	14	233	A
26	14	240	G
26	14	248	G
26	14	249	C
26	14	266	G
26	14	267	C
26	14	269	U
26	14	270(K)	C
26	14	270(L)	U
26	14	270(M)	U
26	14	270(O)	U
26	14	271(B)	G
26	14	271	G
26	14	273(D)	C
26	14	274	G
26	14	275	G
26	14	276	A
26	14	277	C
26	14	278	A
26	14	279	C
26	14	289	A
26	14	290	G
26	14	310	A
26	14	311	A
26	14	319	C
26	14	324	A
26	14	329	G
26	14	330	A
26	14	331	A
26	14	333	G
26	14	346	A

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Mol	Chain	Res	Type
26	14	352	G
26	14	363	G
26	14	363(A)	A
26	14	363(E)	U
26	14	372	G
26	14	382	G
26	14	386	G
26	14	391	G
26	14	395	U
26	14	396	G
26	14	399	G
26	14	405	U
26	14	406	G
26	14	411	G
26	14	412	A
26	14	414	C
26	14	428	A
26	14	443	A
26	14	444	C
26	14	448	U
26	14	454	A
26	14	455	C
26	14	457	A
26	14	459	U
26	14	460	A
26	14	470	A
26	14	471	A
26	14	478	A
26	14	481	G
26	14	483	A
26	14	504	U
26	14	505	A
26	14	509	C
26	14	513	A
26	14	528	A
26	14	529	A
26	14	531	C
26	14	532	A
26	14	533	G
26	14	537	C
26	14	549	G
26	14	556	G

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Mol	Chain	Res	Type
26	14	563	G
26	14	570	G
26	14	573	G
26	14	575	A
26	14	586	A
26	14	587	C
26	14	603	A
26	14	607	U
26	14	609(A)	G
26	14	614	U
26	14	615	G
26	14	617	G
26	14	618	G
26	14	619	G
26	14	620	G
26	14	621	A
26	14	622	G
26	14	627	A
26	14	635	C
26	14	637	A
26	14	645	C
26	14	646	A
26	14	650	C
26	14	651	G
26	14	654	A
26	14	654(E)	C
26	14	654(G)	C
26	14	654(H)	G
26	14	654(I)	C
26	14	654(K)	C
26	14	654(L)	G
26	14	654(T)	A
26	14	656	G
26	14	661	C
26	14	669	G
26	14	686	G
26	14	699	A
26	14	701	G
26	14	717	G
26	14	722	A
26	14	730	C
26	14	738	G

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Mol	Chain	Res	Type
26	14	740	U
26	14	752	A
26	14	753	C
26	14	762	U
26	14	765	G
26	14	775	G
26	14	776	G
26	14	779	U
26	14	782	A
26	14	783	A
26	14	784	A
26	14	785	G
26	14	791	C
26	14	792	G
26	14	800	A
26	14	802	A
26	14	805	G
26	14	812	C
26	14	819	A
26	14	820	A
26	14	827	U
26	14	828	U
26	14	830	G
26	14	832	G
26	14	840	C
26	14	846	C
26	14	848	G
26	14	855	G
26	14	859	G
26	14	865	C
26	14	866	A
26	14	869	G
26	14	878	A
26	14	880	G
26	14	881	G
26	14	882	G
26	14	885	C
26	14	886	C
26	14	887	A
26	14	888	C
26	14	889	C
26	14	890	A

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Mol	Chain	Res	Type
26	14	894	C
26	14	896	A
26	14	897	C
26	14	899	A
26	14	900	A
26	14	901	A
26	14	903	C
26	14	904	C
26	14	906	G
26	14	910	A
26	14	914	C
26	14	915	C
26	14	917	A
26	14	924	C
26	14	932	G
26	14	933	A
26	14	938	G
26	14	941	A
26	14	945	A
26	14	946	G
26	14	953	A
26	14	958	U
26	14	959	A
26	14	961	C
26	14	968	G
26	14	974	G
26	14	980	A
26	14	983	A
26	14	986	C
26	14	989	G
26	14	990	A
26	14	991	C
26	14	996	A
26	14	999	U
26	14	1005	C
26	14	1012	U
26	14	1013	C
26	14	1014	U
26	14	1015	G
26	14	1017	G
26	14	1020	A
26	14	1021	A

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Mol	Chain	Res	Type
26	14	1022	G
26	14	1023	U
26	14	1025	G
26	14	1026	U
26	14	1027	A
26	14	1028	A
26	14	1037	G
26	14	1039	G
26	14	1044	G
26	14	1045	A
26	14	1046	A
26	14	1047	G
26	14	1048	A
26	14	1051	G
26	14	1054	A
26	14	1056	G
26	14	1057	A
26	14	1060	U
26	14	1062	G
26	14	1065	U
26	14	1067	A
26	14	1068	G
26	14	1070	A
26	14	1073	A
26	14	1077	A
26	14	1079	C
26	14	1083	U
26	14	1084	A
26	14	1085	A
26	14	1086	A
26	14	1087	G
26	14	1088	A
26	14	1090	U
26	14	1091	G
26	14	1093	G
26	14	1095	A
26	14	1096	A
26	14	1098	A
26	14	1105	U
26	14	1110	G
26	14	1111	A
26	14	1112	G

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Mol	Chain	Res	Type
26	14	1122	G
26	14	1126	A
26	14	1129	A
26	14	1130	U
26	14	1135	C
26	14	1136	G
26	14	1137	G
26	14	1139	G
26	14	1142	U
26	14	1142(A)	A
26	14	1143	A
26	14	1147	C
26	14	1149	G
26	14	1151	G
26	14	1155	A
26	14	1170	G
26	14	1173	G
26	14	1174	A
26	14	1175	U
26	14	1177	A
26	14	1178	C
26	14	1195	G
26	14	1204	A
26	14	1205	U
26	14	1206	G
26	14	1212	G
26	14	1220	A
26	14	1236	G
26	14	1237	A
26	14	1244	G
26	14	1247	A
26	14	1253	A
26	14	1256	G
26	14	1271	G
26	14	1272	A
26	14	1273	U
26	14	1287	A
26	14	1298	C
26	14	1300	U
26	14	1301	A
26	14	1303	G
26	14	1306	C

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Mol	Chain	Res	Type
26	14	1313	U
26	14	1314	C
26	14	1319	G
26	14	1320	C
26	14	1321	A
26	14	1329	U
26	14	1342	A
26	14	1343	G
26	14	1345	C
26	14	1348	G
26	14	1349	A
26	14	1359	A
26	14	1365	A
26	14	1368	G
26	14	1370	C
26	14	1378	A
26	14	1380	G
26	14	1385	G
26	14	1386	C
26	14	1395	A
26	14	1403	C
26	14	1407	C
26	14	1416	G
26	14	1417	C
26	14	1419	A
26	14	1420	U
26	14	1421	G
26	14	1427	A
26	14	1428	C
26	14	1437	C
26	14	1444(A)	A
26	14	1445	C
26	14	1449	A
26	14	1449(A)	G
26	14	1455	G
26	14	1459	G
26	14	1460	A
26	14	1467	C
26	14	1471	A
26	14	1475	G
26	14	1480	G
26	14	1482	U

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Mol	Chain	Res	Type
26	14	1483	G
26	14	1490	A
26	14	1493	C
26	14	1494	A
26	14	1506	C
26	14	1508	A
26	14	1509	C
26	14	1510	A
26	14	1522	G
26	14	1528	A
26	14	1533	C
26	14	1534	G
26	14	1535	U
26	14	1537	C
26	14	1543	A
26	14	1544	C
26	14	1547	C
26	14	1554	A
26	14	1558	A
26	14	1559	G
26	14	1560	G
26	14	1566	A
26	14	1569	A
26	14	1577	C
26	14	1578	U
26	14	1585	C
26	14	1586	A
26	14	1588	C
26	14	1589	C
26	14	1598	C
26	14	1608	A
26	14	1609	A
26	14	1610	A
26	14	1613	G
26	14	1618	A
26	14	1628	G
26	14	1632	A
26	14	1640	C
26	14	1647	G
26	14	1648	C
26	14	1654	A
26	14	1669	A

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Mol	Chain	Res	Type
26	14	1671	U
26	14	1674	G
26	14	1675	C
26	14	1678	G
26	14	1680	U
26	14	1682	G
26	14	1696	G
26	14	1697	G
26	14	1700	A
26	14	1701	A
26	14	1717	G
26	14	1725	G
26	14	1726	G
26	14	1728	G
26	14	1729	A
26	14	1730	U
26	14	1735	C
26	14	1742	C
26	14	1743	G
26	14	1756	G
26	14	1758	G
26	14	1762	A
26	14	1763	G
26	14	1764	G
26	14	1773	A
26	14	1779	U
26	14	1780	A
26	14	1782	C
26	14	1787	A
26	14	1791	A
26	14	1800	C
26	14	1801	G
26	14	1802	A
26	14	1806	C
26	14	1811	G
26	14	1813	G
26	14	1816	G
26	14	1819	A
26	14	1820	U
26	14	1829	A
26	14	1847	A
26	14	1848	A

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Mol	Chain	Res	Type
26	14	1851	U
26	14	1858	G
26	14	1878	G
26	14	1885	A
26	14	1888	G
26	14	1889	A
26	14	1894	C
26	14	1897	G
26	14	1906	G
26	14	1913	A
26	14	1917	U
26	14	1919	A
26	14	1920	C
26	14	1929	G
26	14	1930	G
26	14	1931	U
26	14	1936	A
26	14	1937	A
26	14	1938	A
26	14	1941	C
26	14	1944	U
26	14	1955	U
26	14	1960	A
26	14	1963	U
26	14	1967	C
26	14	1970	A
26	14	1971	A
26	14	1972	A
26	14	1985	G
26	14	1991	U
26	14	1993	U
26	14	1995	U
26	14	1996	C
26	14	2018	G
26	14	2020	A
26	14	2023	G
26	14	2031	A
26	14	2033	A
26	14	2039	C
26	14	2043	C
26	14	2049	G
26	14	2054	A

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Mol	Chain	Res	Type
26	14	2055	C
26	14	2056	G
26	14	2059	A
26	14	2060	A
26	14	2061	G
26	14	2062	A
26	14	2063	C
26	14	2069	G
26	14	2071	A
26	14	2082	A
26	14	2093	G
26	14	2096	U
26	14	2099	U
26	14	2100	G
26	14	2108	C
26	14	2111	C
26	14	2112	G
26	14	2113	U
26	14	2114	A
26	14	2117	A
26	14	2118	U
26	14	2127	G
26	14	2128	C
26	14	2131	G
26	14	2132	U
26	14	2133	G
26	14	2134	A
26	14	2136	C
26	14	2137	C
26	14	2140	C
26	14	2144	U
26	14	2145	C
26	14	2146	C
26	14	2147	G
26	14	2148	G
26	14	2157	G
26	14	2166	G
26	14	2173	A
26	14	2174	C
26	14	2175	C
26	14	2178	C
26	14	2189	U

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Mol	Chain	Res	Type
26	14	2191	G
26	14	2192	G
26	14	2198	A
26	14	2207	C
26	14	2210	G
26	14	2211	G
26	14	2212	A
26	14	2213	U
26	14	2215	G
26	14	2225	A
26	14	2226	C
26	14	2234	G
26	14	2238	G
26	14	2239	G
26	14	2240	C
26	14	2251	G
26	14	2252	G
26	14	2253	G
26	14	2259	G
26	14	2261	C
26	14	2267	A
26	14	2268	A
26	14	2269	A
26	14	2273	A
26	14	2275	C
26	14	2278	A
26	14	2280	G
26	14	2283	C
26	14	2286	A
26	14	2287	A
26	14	2288	A
26	14	2289	G
26	14	2291	U
26	14	2294	C
26	14	2297	C
26	14	2305	A
26	14	2307	G
26	14	2308	G
26	14	2309	A
26	14	2310	A
26	14	2311	A
26	14	2318	G

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Mol	Chain	Res	Type
26	14	2321	G
26	14	2324	C
26	14	2325	G
26	14	2334	G
26	14	2335	A
26	14	2336	A
26	14	2343	C
26	14	2346	A
26	14	2347	C
26	14	2350	C
26	14	2354	G
26	14	2383	G
26	14	2385	C
26	14	2388	A
26	14	2389	G
26	14	2392	A
26	14	2394	C
26	14	2396	G
26	14	2402	C
26	14	2403	C
26	14	2406	U
26	14	2407	G
26	14	2408	U
26	14	2414	G
26	14	2422	A
26	14	2423	U
26	14	2425	A
26	14	2429	G
26	14	2430	A
26	14	2431	U
26	14	2434	A
26	14	2435	A
26	14	2439	A
26	14	2440	C
26	14	2441	C
26	14	2445	G
26	14	2447	G
26	14	2448	A
26	14	2449	U
26	14	2469	A
26	14	2470	G
26	14	2476	A

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Mol	Chain	Res	Type
26	14	2477	C
26	14	2482	G
26	14	2483	C
26	14	2484	G
26	14	2487	G
26	14	2489	G
26	14	2492	U
26	14	2496	C
26	14	2497	A
26	14	2502	G
26	14	2504	U
26	14	2505	G
26	14	2506	U
26	14	2513	G
26	14	2518	A
26	14	2519	U
26	14	2528	U
26	14	2529	G
26	14	2532	G
26	14	2538	C
26	14	2542	A
26	14	2543	G
26	14	2554	U
26	14	2564	A
26	14	2566	A
26	14	2567	G
26	14	2569	G
26	14	2573	C
26	14	2584	U
26	14	2585	U
26	14	2587	A
26	14	2602	A
26	14	2603	G
26	14	2609	U
26	14	2610	C
26	14	2611	U
26	14	2612	C
26	14	2613	U
26	14	2615	U
26	14	2630	G
26	14	2636	U
26	14	2646	C

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Mol	Chain	Res	Type
26	14	2647	U
26	14	2665	A
26	14	2667	C
26	14	2670	A
26	14	2673	G
26	14	2679	A
26	14	2689	U
26	14	2690	C
26	14	2702	U
26	14	2703	C
26	14	2704	C
26	14	2707	G
26	14	2712(A)	A
26	14	2713	A
26	14	2714	G
26	14	2726	U
26	14	2733	A
26	14	2739	U
26	14	2744	G
26	14	2750	A
26	14	2751	G
26	14	2752	C
26	14	2754	U
26	14	2758	A
26	14	2761	G
26	14	2762	G
26	14	2764	A
26	14	2765	A
26	14	2766	G
26	14	2769	C
26	14	2777	G
26	14	2778	A
26	14	2779	U
26	14	2780	G
26	14	2786	U
26	14	2787	C
26	14	2790	A
26	14	2791	C
26	14	2794	C
26	14	2795	G
26	14	2797	U
26	14	2798	C

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Mol	Chain	Res	Type
26	14	2810	A
26	14	2818	G
26	14	2820	A
26	14	2821	A
26	14	2833	G
26	14	2834	G
26	14	2835	A
26	14	2839	G
26	14	2860	A
26	14	2872	G
26	14	2879	C
26	14	2880	C
26	14	2883	A
26	14	2885	C
26	14	2886	G
26	14	2894	G
26	14	2896	C
26	14	2898	U
26	14	2899	G
27	1J	0	A
27	1J	7	G
27	1J	12	C
27	1J	13	A
27	1J	15	A
27	1J	16	G
27	1J	19	G
27	1J	22	U
27	1J	24	G
27	1J	29	A
27	1J	30	C
27	1J	41	U
27	1J	42	C
27	1J	44	G
27	1J	45	A
27	1J	52	A
27	1J	53	A
27	1J	58	A
27	1J	66	A
27	1J	73	A
27	1J	75	G
27	1J	81	G
27	1J	88	C

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Mol	Chain	Res	Type
27	1J	89	G
27	1J	89(A)	A
27	1J	90	C
27	1J	101	A
27	1J	108	C
27	1J	109	G
27	1J	113	C
27	1J	114	G

All (178) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
1	13	5	U
1	13	50	A
1	13	115	G
1	13	119	A
1	13	244	U
1	13	266	G
1	13	412	A
1	13	422	C
1	13	484	G
1	13	509	A
1	13	560	U
1	13	687	A
1	13	703	G
1	13	748	C
1	13	793	U
1	13	812	C
1	13	991	U
1	13	992	U
1	13	1027	C
1	13	1053	G
1	13	1054	C
1	13	1065	U
1	13	1126	U
1	13	1285	A
1	13	1300	G
1	13	1302	U
1	13	1336	C
1	13	1452	C
1	13	1498	U
1	13	1504	G

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Mol	Chain	Res	Type
22	1K	10	G
22	1K	18	G
23	2K	21	U
23	2K	22	A
23	2K	48	U
24	3K	2	C
24	3K	18	G
24	3K	45	U
24	3K	58	A
26	1H	125	G
26	1H	222	A
26	1H	229	A
26	1H	249	C
26	1H	271(B)	G
26	1H	404	C
26	1H	508	G
26	1H	587	C
26	1H	746	A
26	1H	752	A
26	1H	764	A
26	1H	776	G
26	1H	800	A
26	1H	858	U
26	1H	880	G
26	1H	974(A)	C
26	1H	1022	G
26	1H	1026	U
26	1H	1060	U
26	1H	1085	A
26	1H	1110	G
26	1H	1178	C
26	1H	1301	A
26	1H	1312	U
26	1H	1378	A
26	1H	1396	U
26	1H	1420	U
26	1H	1427	A
26	1H	1508	A
26	1H	1558	A
26	1H	1608	A
26	1H	1609	A
26	1H	1647	G

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Mol	Chain	Res	Type
26	1H	1694	C
26	1H	1699	G
26	1H	1757	U
26	1H	1799	G
26	1H	2060	A
26	1H	2062	A
26	1H	2135	A
26	1H	2157	G
26	1H	2171	A
26	1H	2212	A
26	1H	2225	A
26	1H	2346	A
26	1H	2428	G
26	1H	2439	A
26	1H	2475	C
26	1H	2481	G
26	1H	2566	A
26	1H	2689	U
26	1H	2756	U
1	1G	7	G
1	1G	64	G
1	1G	115	G
1	1G	197	A
1	1G	243	A
1	1G	250	A
1	1G	266	G
1	1G	327	A
1	1G	328	C
1	1G	345	C
1	1G	412	A
1	1G	429	U
1	1G	485	G
1	1G	509	A
1	1G	528	C
1	1G	560	U
1	1G	632	A
1	1G	687	A
1	1G	723	U
1	1G	748	C
1	1G	793	U
1	1G	884	U
1	1G	913	A

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Mol	Chain	Res	Type
1	1G	991	U
1	1G	992	U
1	1G	1053	G
1	1G	1054	C
1	1G	1064	G
1	1G	1126	U
1	1G	1128	C
1	1G	1145	C
1	1G	1157	A
1	1G	1285	A
1	1G	1297	C
1	1G	1300	G
1	1G	1346	A
1	1G	1396	A
1	1G	1453	G
1	1G	1498	U
24	1L	10	G
24	1L	18	G
23	2L	19	G
23	2L	21	U
23	2L	48	U
24	3L	8	U
24	3L	58	A
25	4L	12	A
26	14	34	C
26	14	101	G
26	14	128	C
26	14	196	A
26	14	278	A
26	14	528	A
26	14	685	A
26	14	752	A
26	14	764	A
26	14	774	A
26	14	990	A
26	14	1022	G
26	14	1085	A
26	14	1342	A
26	14	1378	A
26	14	1416	G
26	14	1427	A
26	14	1558	A

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Mol	Chain	Res	Type
26	14	1608	A
26	14	1609	A
26	14	1819	A
26	14	1984	G
26	14	2062	A
26	14	2210	G
26	14	2225	A
26	14	2335	A
26	14	2406	U
26	14	2439	A
26	14	2447	G
26	14	2602	A
26	14	2611	U
26	14	2629	A
26	14	2689	U
26	14	2776	A
26	14	2778	A
26	14	2859	G
26	14	2893	G
27	1J	12	C
27	1J	56	G
27	1J	88	C

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

17 non-standard protein/DNA/RNA residues are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
22	PSU	1K	32	22	15,21,22	1.01	1 (6%)	16,30,33	2.05	2 (12%)
22	MIA	1K	37	22	22,31,32	0.89	1 (4%)	26,44,47	1.73	6 (23%)
22	PSU	1K	39	22	15,21,22	0.98	1 (6%)	16,30,33	2.02	4 (25%)
22	7MG	1K	46	22	20,26,27	3.26	6 (30%)	23,39,42	2.23	5 (21%)
22	5MU	1K	54	22	13,22,23	1.70	2 (15%)	16,32,35	1.54	1 (6%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
22	PSU	1K	55	22	15,21,22	1.19	1 (6%)	16,30,33	2.30	3 (18%)
22	4SU	1K	8	22	12,21,22	3.26	2 (16%)	15,30,33	1.07	1 (6%)
23	OMC	2K	33	23	15,22,23	2.16	4 (26%)	20,31,34	1.58	2 (10%)
23	7MG	2K	47	23	20,26,27	3.30	6 (30%)	23,39,42	2.28	9 (39%)
23	5MU	2K	55	23	13,22,23	1.62	2 (15%)	16,32,35	1.29	1 (6%)
23	PSU	2K	56	23	15,21,22	0.98	1 (6%)	16,30,33	1.76	4 (25%)
23	4SU	2K	8	23	12,21,22	3.11	2 (16%)	15,30,33	1.02	1 (6%)
23	OMC	2L	33	23	15,22,23	2.23	4 (26%)	20,31,34	2.16	4 (20%)
23	7MG	2L	47	23	20,26,27	3.34	5 (25%)	23,39,42	2.12	5 (21%)
23	5MU	2L	55	23	13,22,23	1.68	2 (15%)	16,32,35	1.18	1 (6%)
23	PSU	2L	56	23	15,21,22	1.11	1 (6%)	16,30,33	1.85	3 (18%)
23	4SU	2L	8	23	12,21,22	3.32	2 (16%)	15,30,33	1.23	1 (6%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
22	PSU	1K	32	22	-	0/7/25/26	0/2/2/2
22	MIA	1K	37	22	-	0/11/33/34	0/3/3/3
22	PSU	1K	39	22	-	0/7/25/26	0/2/2/2
22	7MG	1K	46	22	-	0/7/37/38	0/3/3/3
22	5MU	1K	54	22	-	0/3/25/26	0/2/2/2
22	PSU	1K	55	22	-	0/7/25/26	0/2/2/2
22	4SU	1K	8	22	-	0/3/25/26	0/2/2/2
23	OMC	2K	33	23	-	0/5/27/28	0/2/2/2
23	7MG	2K	47	23	-	0/7/37/38	0/3/3/3
23	5MU	2K	55	23	-	0/3/25/26	0/2/2/2
23	PSU	2K	56	23	-	0/7/25/26	0/2/2/2
23	4SU	2K	8	23	-	0/3/25/26	0/2/2/2
23	OMC	2L	33	23	-	0/5/27/28	0/2/2/2
23	7MG	2L	47	23	-	0/7/37/38	0/3/3/3
23	5MU	2L	55	23	-	0/3/25/26	0/2/2/2
23	PSU	2L	56	23	-	0/7/25/26	0/2/2/2
23	4SU	2L	8	23	-	0/3/25/26	0/2/2/2

All (43) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
23	2K	47	7MG	C5-C4	-5.99	1.23	1.39
23	2L	47	7MG	C5-C4	-5.75	1.23	1.39
22	1K	46	7MG	C5-C4	-5.32	1.25	1.39
23	2K	55	5MU	C4-N3	-3.15	1.27	1.33
23	2L	55	5MU	C4-N3	-3.15	1.27	1.33
22	1K	54	5MU	C4-N3	-2.82	1.27	1.33
22	1K	37	MIA	C4-N3	-2.19	1.32	1.35
22	1K	46	7MG	C1'-N9	2.05	1.49	1.44
23	2K	47	7MG	C2-N1	2.44	1.40	1.35
23	2K	56	PSU	C4-N3	2.75	1.38	1.33
22	1K	39	PSU	C4-N3	2.94	1.38	1.33
22	1K	32	PSU	C4-N3	3.01	1.38	1.33
23	2L	33	OMC	C4-N4	3.25	1.44	1.35
22	1K	46	7MG	C2-N2	3.37	1.41	1.34
23	2K	33	OMC	C4-N4	3.40	1.44	1.35
23	2K	33	OMC	C2-N3	3.42	1.45	1.38
23	2L	56	PSU	C4-N3	3.45	1.39	1.33
23	2K	47	7MG	C2-N2	3.53	1.41	1.34
23	2L	47	7MG	C2-N2	3.55	1.41	1.34
22	1K	55	PSU	C4-N3	3.72	1.39	1.33
23	2L	33	OMC	C2-N3	3.95	1.46	1.38
23	2L	33	OMC	C5-C4	4.02	1.50	1.41
23	2K	33	OMC	C5-C4	4.40	1.51	1.41
23	2K	55	5MU	C2-N3	4.44	1.47	1.38
22	1K	46	7MG	C8-N7	4.47	1.64	1.43
23	2L	47	7MG	C8-N7	4.57	1.64	1.43
23	2K	47	7MG	C8-N7	4.64	1.65	1.43
23	2L	55	5MU	C2-N3	4.77	1.48	1.38
23	2K	47	7MG	C6-C5	4.82	1.48	1.41
23	2K	33	OMC	C6-N1	4.86	1.42	1.35
22	1K	54	5MU	C2-N3	5.03	1.48	1.38
22	1K	46	7MG	C6-C5	5.16	1.48	1.41
23	2L	47	7MG	C6-C5	5.32	1.49	1.41
23	2L	33	OMC	C6-N1	5.40	1.42	1.35
23	2K	8	4SU	C6-N1	5.41	1.42	1.35
23	2L	8	4SU	C6-N1	7.00	1.44	1.35
22	1K	8	4SU	C6-N1	7.00	1.44	1.35
22	1K	8	4SU	C5-C4	8.63	1.49	1.38
23	2L	8	4SU	C5-C4	8.89	1.50	1.38
23	2K	8	4SU	C5-C4	9.03	1.50	1.38
23	2K	47	7MG	C4-N3	10.37	1.47	1.34
22	1K	46	7MG	C4-N3	10.65	1.47	1.34
23	2L	47	7MG	C4-N3	10.84	1.48	1.34

All (53) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
23	2L	47	7MG	C5-C4-N3	-6.34	120.29	126.74
22	1K	46	7MG	C5-C4-N3	-6.25	120.38	126.74
22	1K	37	MIA	C5-C6-N1	-4.81	115.70	120.58
23	2L	8	4SU	C5-C4-N3	-4.25	119.05	123.56
23	2K	47	7MG	C5-C4-N3	-3.90	122.77	126.74
22	1K	55	PSU	C5-C1'-C2'	-3.71	109.13	115.44
23	2K	47	7MG	C4-N9-C1'	-3.63	118.04	126.65
22	1K	8	4SU	C5-C4-N3	-3.49	119.86	123.56
23	2K	47	7MG	N1-C2-N3	-3.39	119.97	125.51
23	2K	47	7MG	C5-C6-N1	-3.24	118.56	123.39
23	2L	47	7MG	N1-C2-N3	-3.24	120.22	125.51
23	2K	47	7MG	N3-C4-N9	-3.22	122.82	126.98
23	2L	56	PSU	C5-C6-N1	-3.19	119.92	124.38
23	2K	8	4SU	C5-C4-N3	-3.06	120.32	123.56
22	1K	46	7MG	N1-C2-N3	-3.04	120.55	125.51
22	1K	46	7MG	C5-C6-N1	-2.94	119.01	123.39
23	2L	33	OMC	C5-C4-N3	-2.71	118.36	121.79
23	2L	47	7MG	C5-C6-N1	-2.65	119.45	123.39
23	2K	56	PSU	C5-C6-N1	-2.42	121.01	124.38
22	1K	39	PSU	C5-C1'-C2'	-2.38	111.39	115.44
23	2L	33	OMC	C6-N1-C2	-2.21	117.73	121.33
22	1K	39	PSU	C5-C6-N1	-2.16	121.37	124.38
23	2K	56	PSU	C5-C1'-C2'	2.01	118.84	115.44
23	2K	56	PSU	O4'-C1'-C2'	2.04	106.89	104.69
22	1K	37	MIA	N6-C6-N1	2.05	120.99	118.55
22	1K	32	PSU	O4'-C1'-C2'	2.05	106.91	104.69
23	2K	47	7MG	C8-N9-C1'	2.07	128.63	122.43
22	1K	55	PSU	O4'-C1'-C2'	2.13	106.99	104.69
23	2L	56	PSU	O4'-C1'-C2'	2.15	107.02	104.69
22	1K	37	MIA	C1'-N9-C4	2.39	129.47	126.81
23	2K	33	OMC	N4-C4-N3	2.63	121.10	116.50
22	1K	39	PSU	O4'-C1'-C2'	2.79	107.70	104.69
23	2K	47	7MG	N2-C2-N1	2.91	122.00	117.20
22	1K	37	MIA	C2-N1-C6	2.96	121.26	113.13
22	1K	37	MIA	C12-N6-C6	3.23	127.19	123.46
23	2L	47	7MG	C6-N1-C2	3.30	119.75	115.88
22	1K	46	7MG	C6-N1-C2	3.62	120.13	115.88
22	1K	37	MIA	C13-C12-N6	3.66	119.09	112.25
23	2L	33	OMC	N4-C4-N3	3.73	123.02	116.50
23	2K	47	7MG	C6-N1-C2	4.05	120.63	115.88
23	2L	55	5MU	C4-N3-C2	4.07	118.55	115.16
23	2L	47	7MG	C5-C4-N9	4.51	113.53	106.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
23	2K	55	5MU	C4-N3-C2	4.56	118.96	115.16
22	1K	46	7MG	C5-C4-N9	5.01	114.33	106.25
23	2K	47	7MG	C5-C4-N9	5.05	114.41	106.25
23	2K	56	PSU	C4-N3-C2	5.19	119.49	115.16
23	2L	56	PSU	C4-N3-C2	5.34	119.61	115.16
23	2K	33	OMC	C6-C5-C4	5.55	119.61	117.44
22	1K	54	5MU	C4-N3-C2	5.65	119.87	115.16
22	1K	39	PSU	C4-N3-C2	6.16	120.30	115.16
22	1K	32	PSU	C4-N3-C2	7.23	121.19	115.16
22	1K	55	PSU	C4-N3-C2	7.48	121.40	115.16
23	2L	33	OMC	C6-C5-C4	7.79	120.49	117.44

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

12 monomers are involved in 25 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
22	1K	37	MIA	3	0
22	1K	46	7MG	3	0
22	1K	54	5MU	1	0
22	1K	55	PSU	1	0
23	2K	33	OMC	1	0
23	2K	47	7MG	2	0
23	2K	55	5MU	3	0
23	2K	8	4SU	2	0
23	2L	33	OMC	5	0
23	2L	47	7MG	1	0
23	2L	55	5MU	3	0
23	2L	8	4SU	1	0

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 1253 ligands modelled in this entry, 1253 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å²)	Q<0.9
1	13	1498/1522 (98%)	0.21	44 (2%)	55	29	73, 122, 204, 309	0
1	1G	1497/1522 (98%)	0.24	72 (4%)	34	15	92, 145, 218, 318	0
2	12	237/256 (92%)	0.20	17 (7%)	18	6	168, 199, 220, 228	0
2	1E	237/256 (92%)	0.12	8 (3%)	49	23	132, 167, 193, 205	0
3	22	206/239 (86%)	0.04	7 (3%)	49	23	161, 182, 205, 218	0
3	2E	205/239 (85%)	-0.02	1 (0%)	91	81	103, 128, 161, 170	0
4	32	208/209 (99%)	0.59	18 (8%)	13	4	122, 144, 165, 174	0
4	3E	208/209 (99%)	0.78	31 (14%)	3	1	102, 129, 151, 165	0
5	42	151/162 (93%)	0.52	18 (11%)	6	2	136, 154, 171, 203	0
5	4E	151/162 (93%)	0.57	19 (12%)	5	2	97, 120, 143, 176	0
6	52	101/101 (100%)	-0.27	0	100	100	106, 127, 147, 161	0
6	5E	101/101 (100%)	-0.07	0	100	100	101, 123, 148, 161	0
7	62	155/156 (99%)	0.92	27 (17%)	2	1	136, 154, 182, 207	0
7	6E	155/156 (99%)	0.49	16 (10%)	9	3	124, 139, 168, 193	0
8	72	138/138 (100%)	1.27	41 (29%)	1	0	130, 158, 174, 179	0
8	7E	138/138 (100%)	0.89	26 (18%)	2	0	105, 129, 141, 149	0
9	82	127/128 (99%)	3.26	75 (59%)	0	0	141, 183, 204, 208	0
9	8E	127/128 (99%)	1.66	45 (35%)	0	0	107, 156, 179, 193	0
10	1A	99/105 (94%)	1.96	39 (39%)	0	0	153, 184, 203, 210	0
10	1I	99/105 (94%)	1.40	32 (32%)	1	0	102, 155, 187, 192	0
11	2A	116/129 (89%)	0.51	11 (9%)	10	4	111, 137, 158, 182	0
11	2I	116/129 (89%)	0.27	8 (6%)	20	7	90, 127, 153, 180	0
12	3A	125/132 (94%)	1.25	35 (28%)	1	0	109, 134, 160, 185	0
12	3I	125/132 (94%)	0.45	11 (8%)	12	4	83, 96, 131, 179	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
13	4A	117/126 (92%)	1.15	30 (25%) 1 0	140, 184, 206, 216	0
13	4I	118/126 (93%)	0.25	7 (5%) 26 10	101, 139, 159, 170	0
14	5A	58/61 (95%)	4.96	46 (79%) 0 0	164, 179, 194, 201	0
14	5I	60/61 (98%)	1.88	23 (38%) 0 0	106, 118, 137, 147	0
15	6A	88/89 (98%)	0.81	12 (13%) 4 1	118, 139, 155, 159	0
15	6I	88/89 (98%)	1.01	21 (23%) 1 0	98, 125, 142, 154	0
16	7A	84/88 (95%)	1.03	19 (22%) 1 0	113, 132, 150, 174	0
16	7I	84/88 (95%)	3.10	55 (65%) 0 0	122, 135, 169, 187	0
17	8A	100/105 (95%)	1.17	21 (21%) 1 0	118, 138, 153, 176	0
17	8I	100/105 (95%)	0.66	14 (14%) 4 1	108, 127, 140, 145	0
18	9A	72/88 (81%)	0.05	2 (2%) 56 30	116, 142, 178, 198	0
18	9I	72/88 (81%)	0.10	0 100 100	109, 127, 162, 187	0
19	AA	78/93 (83%)	1.25	25 (32%) 1 0	172, 203, 216, 222	0
19	AI	81/93 (87%)	0.47	4 (4%) 33 14	109, 135, 160, 166	0
20	BA	99/106 (93%)	1.63	43 (43%) 0 0	108, 130, 158, 173	0
20	BI	99/106 (93%)	1.35	34 (34%) 0 0	129, 144, 177, 183	0
21	1B	25/27 (92%)	6.09	23 (92%) 0 0	147, 163, 179, 196	0
21	1F	25/27 (92%)	4.76	22 (88%) 0 0	114, 126, 139, 164	0
22	1K	69/76 (90%)	0.13	5 (7%) 18 6	101, 222, 265, 271	0
23	2K	72/77 (93%)	0.05	0 100 100	84, 108, 137, 146	0
23	2L	72/77 (93%)	-0.41	1 (1%) 78 57	96, 136, 166, 170	0
24	1L	76/76 (100%)	0.82	13 (17%) 2 1	147, 264, 289, 295	0
24	3K	76/76 (100%)	-0.06	2 (2%) 59 33	92, 255, 286, 289	0
24	3L	76/76 (100%)	0.50	9 (11%) 6 2	104, 264, 294, 299	0
25	4K	13/27 (48%)	0.92	0 100 100	88, 101, 151, 157	0
25	4L	9/27 (33%)	0.60	0 100 100	121, 154, 165, 175	0
26	14	2909/2917 (99%)	0.21	70 (2%) 62 37	67, 105, 270, 378	0
26	1H	2912/2917 (99%)	0.28	12 (0%) 93 84	56, 91, 253, 353	0
27	16	122/122 (100%)	-0.17	0 100 100	83, 111, 133, 217	0
27	1J	122/122 (100%)	-0.14	1 (0%) 87 72	107, 152, 174, 226	0
28	11	272/276 (98%)	0.50	8 (2%) 55 29	57, 83, 100, 110	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
28	19	273/276 (98%)	0.91	41 (15%) 3 1	67, 94, 110, 125	0
29	21	205/206 (99%)	0.90	27 (13%) 4 1	66, 109, 151, 165	0
29	29	205/206 (99%)	1.29	67 (32%) 1 0	76, 113, 159, 183	0
30	31	202/210 (96%)	0.32	2 (0%) 84 66	60, 94, 130, 150	0
30	39	208/210 (99%)	0.43	14 (6%) 21 7	76, 123, 179, 199	0
31	41	181/182 (99%)	0.30	9 (4%) 32 13	100, 121, 159, 172	0
31	49	181/182 (99%)	1.07	41 (22%) 1 0	148, 169, 198, 209	0
32	51	174/180 (96%)	0.12	5 (2%) 55 29	99, 122, 139, 150	0
32	59	169/180 (93%)	1.63	62 (36%) 0 0	169, 218, 243, 258	0
33	61	146/148 (98%)	-0.02	4 (2%) 58 32	98, 155, 170, 175	0
33	69	146/148 (98%)	0.71	22 (15%) 3 1	102, 146, 170, 175	0
34	15	138/140 (98%)	1.27	35 (25%) 1 0	98, 129, 161, 185	0
34	58	138/140 (98%)	0.79	15 (10%) 7 2	81, 109, 148, 163	0
35	25	122/122 (100%)	0.51	9 (7%) 17 6	85, 107, 125, 135	0
35	68	122/122 (100%)	0.41	3 (2%) 61 35	75, 92, 111, 124	0
36	35	150/150 (100%)	1.13	40 (26%) 1 0	77, 128, 162, 198	0
36	78	150/150 (100%)	0.56	11 (7%) 18 6	64, 98, 124, 176	0
37	45	141/141 (100%)	2.38	78 (55%) 0 0	97, 129, 154, 167	0
37	88	138/141 (97%)	0.52	7 (5%) 32 13	70, 95, 115, 147	0
38	55	117/118 (99%)	0.97	20 (17%) 2 1	77, 97, 113, 133	0
38	98	118/118 (100%)	0.82	13 (11%) 7 2	79, 102, 125, 139	0
39	65	111/112 (99%)	1.38	36 (32%) 1 0	118, 145, 162, 172	0
39	A8	111/112 (99%)	0.35	4 (3%) 46 21	92, 107, 130, 144	0
40	75	137/146 (93%)	0.31	10 (7%) 18 6	96, 115, 173, 208	0
40	B8	137/146 (93%)	0.49	12 (8%) 12 4	86, 111, 170, 202	0
41	85	117/118 (99%)	0.79	17 (14%) 3 1	86, 115, 155, 180	0
41	C8	117/118 (99%)	0.87	19 (16%) 3 1	71, 98, 136, 147	0
42	95	101/101 (100%)	0.51	14 (13%) 4 1	86, 143, 161, 177	0
42	D8	101/101 (100%)	0.31	10 (9%) 9 3	73, 123, 149, 162	0
43	A5	113/113 (100%)	0.42	7 (6%) 24 9	77, 94, 124, 183	0
43	E8	113/113 (100%)	0.38	3 (2%) 58 32	75, 91, 125, 164	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
44	B5	92/96 (95%)	0.68	8 (8%) 13 4	85, 102, 129, 133	0
44	F8	94/96 (97%)	0.29	1 (1%) 82 63	69, 88, 113, 128	0
45	C5	104/110 (94%)	1.29	25 (24%) 1 0	107, 137, 167, 175	0
45	G8	104/110 (94%)	0.30	0 100 100	87, 113, 147, 166	0
46	D5	179/206 (86%)	0.99	35 (19%) 1 0	139, 176, 263, 269	0
46	H8	175/206 (84%)	-0.05	1 (0%) 90 78	97, 139, 228, 236	0
47	E5	76/85 (89%)	1.89	35 (46%) 0 0	85, 111, 130, 169	0
47	I8	80/85 (94%)	1.09	19 (23%) 1 0	71, 90, 122, 131	0
48	F5	97/98 (98%)	1.74	32 (32%) 0 0	79, 102, 146, 163	0
48	J8	97/98 (98%)	1.14	14 (14%) 3 1	69, 91, 145, 176	0
49	G5	66/72 (91%)	0.19	2 (3%) 54 27	101, 121, 141, 167	0
49	K8	67/72 (93%)	0.39	2 (2%) 54 27	76, 96, 119, 154	0
50	H5	59/60 (98%)	1.04	13 (22%) 1 0	97, 126, 167, 182	0
50	L8	57/60 (95%)	0.18	0 100 100	77, 98, 120, 133	0
51	I5	63/71 (88%)	1.49	22 (34%) 0 0	179, 219, 236, 243	0
51	M8	66/71 (92%)	0.44	6 (9%) 11 4	128, 174, 203, 214	0
52	J5	56/60 (93%)	0.38	4 (7%) 19 6	77, 103, 150, 162	0
52	N8	58/60 (96%)	0.93	6 (10%) 9 3	68, 115, 186, 191	0
53	K5	45/54 (83%)	7.08	38 (84%) 0 0	153, 180, 196, 202	0
53	O8	45/54 (83%)	3.61	32 (71%) 0 0	131, 159, 177, 183	0
54	L5	46/49 (93%)	0.72	3 (6%) 22 8	64, 76, 89, 102	0
54	P8	45/49 (91%)	0.14	0 100 100	59, 64, 78, 92	0
55	M5	60/65 (92%)	2.19	37 (61%) 0 0	89, 101, 125, 147	0
55	Q8	61/65 (93%)	1.77	23 (37%) 0 0	74, 91, 117, 128	0
All	All	21042/21688 (97%)	0.57	2138 (10%) 9 3	56, 122, 214, 378	0

All (2138) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
53	K5	42	TRP	19.7
53	K5	51	GLU	19.1
24	3L	17	C	18.5
7	62	81	GLY	16.9
7	62	82	GLY	16.6

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Mol	Chain	Res	Type	RSRZ
53	K5	13	CYS	16.3
39	65	2	ALA	16.2
24	1L	76	A	13.2
8	72	1	MET	13.1
13	4A	102	ARG	12.5
53	K5	12	GLU	12.5
9	82	115	GLY	12.4
48	F5	98	LEU	12.4
53	K5	50	ARG	11.9
14	5A	10	ALA	11.7
53	K5	53	LYS	11.7
46	D5	179	ASP	11.6
53	K5	14	THR	11.6
14	5A	34	TYR	11.4
48	F5	97	LEU	11.4
48	J8	98	LEU	11.2
37	45	1	MET	11.2
21	1B	14	TRP	10.8
53	K5	22	ALA	10.7
14	5A	31	ARG	10.7
24	1L	71	G	10.6
10	1A	59	SER	10.5
10	1A	46	ARG	10.5
7	6E	81	GLY	10.5
21	1F	26	LYS	10.4
53	K5	41	PRO	10.3
14	5A	38	GLY	10.3
21	1F	17	THR	10.3
14	5A	39	LEU	10.2
21	1B	13	ILE	10.1
9	82	123	PRO	10.1
9	82	110	GLU	10.0
53	K5	21	TYR	10.0
9	82	66	ARG	9.9
14	5A	32	SER	9.8
10	1A	47	PHE	9.7
53	K5	49	HIS	9.7
53	K5	24	GLU	9.7
53	K5	9	LEU	9.6
53	K5	39	TYR	9.6
16	7I	32	TYR	9.5
7	62	80	VAL	9.4

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Mol	Chain	Res	Type	RSRZ
53	K5	26	ASN	9.4
9	82	69	GLY	9.3
52	N8	59	GLU	9.3
9	82	109	VAL	9.1
32	59	169	VAL	9.1
24	3L	34	G	9.1
9	82	116	LYS	9.0
53	O8	42	TRP	8.9
53	K5	52	VAL	8.8
19	AA	79	THR	8.8
21	1B	26	LYS	8.8
9	82	126	SER	8.7
32	59	96	ALA	8.7
53	K5	25	LYS	8.7
21	1B	25	LYS	8.7
16	7I	22	THR	8.7
9	82	127	LYS	8.6
21	1B	2	GLY	8.6
24	1L	72	C	8.6
32	59	4	ILE	8.6
10	1A	58	ASP	8.5
51	M8	55	ARG	8.5
53	K5	23	THR	8.5
16	7I	7	ALA	8.5
9	8E	126	SER	8.5
14	5A	35	ARG	8.4
19	AA	78	ARG	8.4
10	1A	55	LYS	8.4
55	Q8	34	TRP	8.3
14	5A	59	ALA	8.3
46	D5	178	GLU	8.3
9	82	114	TYR	8.3
10	1A	54	PHE	8.2
21	1B	24	ARG	8.2
9	82	12	GLU	8.2
11	2A	11	LYS	8.2
26	14	2901	C	8.1
14	5A	36	PHE	8.1
14	5A	37	PHE	8.1
13	4A	101	GLN	8.1
21	1F	16	GLY	8.1
33	69	1	MET	8.1

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Mol	Chain	Res	Type	RSRZ
39	A8	2	ALA	8.0
21	1B	9	ARG	8.0
20	BI	70	SER	8.0
17	8A	101	ARG	8.0
9	82	125	TYR	7.9
13	4A	6	GLY	7.9
21	1B	6	ARG	7.9
26	14	2902	C	7.9
36	35	150	ALA	7.9
32	59	94	TYR	7.8
46	D5	121	HIS	7.8
35	25	1	MET	7.8
9	82	128	ARG	7.8
26	14	2146	C	7.8
53	O8	49	HIS	7.8
9	82	14	VAL	7.8
37	45	91	GLU	7.7
53	K5	40	CYS	7.7
22	1K	17	C	7.6
53	O8	25	LYS	7.6
14	5A	58	LYS	7.5
21	1F	15	ARG	7.5
52	N8	60	VAL	7.5
9	8E	111	ARG	7.5
20	BI	72	LEU	7.4
21	1B	21	TYR	7.4
29	29	150	VAL	7.4
26	14	2147	G	7.4
21	1F	2	GLY	7.4
22	1K	76	A	7.4
10	1A	64	GLU	7.3
53	K5	20	ASN	7.3
9	82	71	SER	7.3
9	82	70	LYS	7.3
14	5A	41	ARG	7.3
21	1B	22	ARG	7.3
16	7I	31	LYS	7.2
9	82	65	VAL	7.2
32	59	100	GLY	7.2
37	45	104	PHE	7.2
14	5A	33	VAL	7.1
8	7E	1	MET	7.1

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Mol	Chain	Res	Type	RSRZ
48	J8	97	LEU	7.1
24	1L	75	C	7.1
21	1B	16	GLY	7.1
53	O8	18	ARG	7.1
37	45	65	PHE	7.1
29	29	151	TYR	7.1
53	O8	53	LYS	7.0
53	K5	19	ARG	7.0
8	72	2	LEU	7.0
54	L5	1	MET	6.9
16	7I	29	ASP	6.9
45	C5	59	GLY	6.9
9	82	106	ALA	6.9
14	5A	61	TRP	6.9
37	45	103	MET	6.9
9	8E	110	GLU	6.8
53	O8	20	ASN	6.8
1	1G	1286	A	6.8
10	1I	5	ARG	6.8
10	1I	64	GLU	6.8
21	1F	14	TRP	6.8
13	4A	103	THR	6.8
52	N8	58	LEU	6.8
53	K5	18	ARG	6.7
53	K5	11	LEU	6.7
48	J8	92	LYS	6.7
32	59	132	ARG	6.7
51	M8	66	SER	6.7
9	82	117	HIS	6.7
14	5A	30	ALA	6.7
39	65	3	ARG	6.6
10	1A	69	ASN	6.6
21	1B	10	ARG	6.6
47	E5	75	LEU	6.6
21	1F	3	LYS	6.6
32	59	95	ARG	6.6
9	82	64	THR	6.6
36	35	110	TYR	6.5
16	7I	9	PHE	6.5
20	BI	18	GLN	6.5
10	1A	62	HIS	6.5
39	65	5	THR	6.5

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Mol	Chain	Res	Type	RSRZ
12	3A	19	ARG	6.5
9	8E	36	TYR	6.5
9	8E	115	GLY	6.5
11	2A	13	GLN	6.5
14	5A	9	LYS	6.5
16	7I	59	TRP	6.5
32	59	115	VAL	6.5
31	49	138	GLN	6.5
21	1F	6	ARG	6.5
38	55	5	LYS	6.4
14	5A	21	TYR	6.4
14	5A	4	LYS	6.4
32	59	107	VAL	6.4
53	O8	44	ARG	6.4
13	4A	4	ILE	6.4
31	49	34	LEU	6.3
10	1A	40	LEU	6.3
9	8E	117	HIS	6.3
33	69	36	ALA	6.3
46	D5	146	ILE	6.3
9	82	113	LYS	6.3
7	62	79	ARG	6.3
9	82	36	TYR	6.3
9	8E	127	LYS	6.3
29	21	78	LEU	6.3
9	82	10	ARG	6.3
32	59	155	SER	6.3
9	82	111	ARG	6.2
48	F5	96	LYS	6.2
16	7I	28	ARG	6.2
51	I5	63	TYR	6.2
37	45	17	LEU	6.2
14	5A	25	VAL	6.2
12	3A	28	LYS	6.2
20	BA	69	GLY	6.2
7	62	41	ARG	6.2
34	15	75	TYR	6.2
47	E5	76	GLY	6.2
9	82	118	LYS	6.1
53	K5	10	LEU	6.1
24	1L	73	A	6.1
21	1B	23	PRO	6.1

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Mol	Chain	Res	Type	RSRZ
21	1B	18	TYR	6.1
11	2I	11	LYS	6.1
55	M5	40	GLU	6.1
13	4A	97	PRO	6.1
14	5A	22	THR	6.1
37	45	38	GLU	6.1
16	7I	27	LYS	6.0
7	6E	83	ALA	6.0
10	1A	48	THR	6.0
14	5A	23	ARG	6.0
33	69	38	LEU	6.0
24	3K	17	C	6.0
12	3A	20	LYS	6.0
21	1F	18	TYR	6.0
10	1A	65	LEU	6.0
12	3I	19	ARG	6.0
24	1L	17	C	6.0
46	D5	115	GLY	5.9
53	K5	36	LEU	5.9
14	5A	29	ARG	5.9
16	7I	6	LEU	5.9
26	1H	2477	C	5.9
20	BA	70	SER	5.9
53	O8	23	THR	5.9
21	1B	15	ARG	5.9
32	59	89	ILE	5.9
53	K5	45	LYS	5.9
37	45	66	ILE	5.9
9	82	112	LYS	5.9
2	1E	14	GLY	5.8
10	1I	66	ARG	5.8
40	B8	106	SER	5.8
9	8E	8	GLY	5.8
19	AA	82	GLY	5.8
9	82	13	ALA	5.8
37	45	10	ARG	5.8
53	O8	9	LEU	5.8
9	82	7	THR	5.7
13	4A	99	ARG	5.7
16	7I	1	MET	5.7
14	5A	55	GLY	5.7
45	C5	46	LYS	5.7

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Mol	Chain	Res	Type	RSRZ
14	5A	19	ARG	5.7
8	7E	2	LEU	5.7
48	J8	93	GLU	5.7
13	4A	66	LEU	5.7
9	82	108	VAL	5.7
29	29	205	ALA	5.7
29	29	117	MET	5.7
39	65	33	LYS	5.6
9	82	120	ARG	5.6
10	1I	46	ARG	5.6
37	45	68	ILE	5.6
7	6E	82	GLY	5.6
16	7I	66	PRO	5.6
10	1I	60	ARG	5.6
16	7I	62	VAL	5.6
33	69	35	LEU	5.6
53	K5	48	VAL	5.6
11	2A	12	ARG	5.6
16	7I	18	ARG	5.6
19	AA	80	TYR	5.6
53	O8	13	CYS	5.6
47	E5	21	LEU	5.6
29	21	205	ALA	5.5
16	7I	65	GLN	5.5
9	82	119	ALA	5.5
34	15	84	LYS	5.5
33	69	3	VAL	5.5
9	82	121	ARG	5.5
40	B8	1	MET	5.5
12	3A	64	TYR	5.5
1	13	1286	A	5.5
32	59	103	LEU	5.5
7	6E	79	ARG	5.5
27	1J	88	C	5.5
32	59	90	LYS	5.5
26	1H	2799	A	5.4
14	5A	26	ARG	5.4
1	1G	973	G	5.4
22	1K	73	A	5.4
7	62	37	ASN	5.4
14	5A	8	GLU	5.4
51	I5	52	THR	5.4

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Mol	Chain	Res	Type	RSRZ
14	5A	60	SER	5.4
31	49	90	LEU	5.4
38	55	4	LEU	5.4
9	82	37	PHE	5.3
8	7E	3	THR	5.3
8	72	3	THR	5.3
9	8E	114	TYR	5.3
10	1A	60	ARG	5.3
45	C5	47	LYS	5.3
47	E5	71	ASP	5.3
14	5A	11	LYS	5.3
36	35	65	ARG	5.3
21	1B	5	ASP	5.3
2	12	132	LYS	5.3
16	7I	12	LYS	5.3
10	1A	57	LYS	5.2
21	1F	22	ARG	5.2
12	3A	32	PHE	5.2
42	95	74	LYS	5.2
8	72	89	PRO	5.2
37	45	61	GLY	5.2
10	1I	62	HIS	5.2
17	8I	36	ILE	5.2
14	5I	61	TRP	5.2
9	82	8	GLY	5.2
7	62	78	ARG	5.2
16	7I	30	GLY	5.2
16	7I	17	TYR	5.2
47	I8	43	THR	5.2
10	1A	56	HIS	5.2
9	8E	15	ALA	5.2
26	14	1092	C	5.2
37	45	90	VAL	5.2
13	4A	100	GLY	5.1
53	K5	34	LEU	5.1
38	55	9	LYS	5.1
8	72	133	LEU	5.1
20	BI	15	ARG	5.1
34	15	85	ILE	5.1
34	58	72	TYR	5.1
45	C5	45	VAL	5.1
9	82	67	GLY	5.1

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Mol	Chain	Res	Type	RSRZ
7	62	83	ALA	5.1
9	82	42	ARG	5.1
29	29	114	ALA	5.1
10	1I	65	LEU	5.0
53	K5	37	ARG	5.0
33	69	2	LYS	5.0
37	45	34	LEU	5.0
53	K5	46	HIS	5.0
33	69	4	ILE	5.0
48	J8	96	LYS	5.0
46	D5	119	GLU	5.0
46	D5	117	LEU	5.0
32	59	164	TYR	5.0
10	1A	67	THR	5.0
9	8E	106	ALA	5.0
21	1F	25	LYS	5.0
32	59	88	LEU	5.0
14	5A	17	LYS	5.0
36	35	18	ARG	4.9
13	4I	102	ARG	4.9
55	M5	16	ILE	4.9
21	1B	11	GLY	4.9
34	15	80	GLY	4.9
16	7I	39	TYR	4.9
34	15	108	PRO	4.9
12	3A	21	LYS	4.9
32	59	92	ILE	4.9
2	12	133	LYS	4.9
26	14	2119	A	4.9
9	8E	116	LYS	4.9
34	58	83	LYS	4.9
21	1B	8	THR	4.9
53	O8	26	ASN	4.9
32	59	99	VAL	4.9
12	3A	128	ALA	4.9
53	O8	14	THR	4.9
53	K5	16	CYS	4.9
10	1A	39	PRO	4.9
55	M5	22	VAL	4.9
17	8I	26	GLN	4.8
9	82	15	ALA	4.8
31	49	35	GLU	4.8

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Mol	Chain	Res	Type	RSRZ
31	49	75	LYS	4.8
16	7I	4	ILE	4.8
16	7I	19	ILE	4.8
15	6A	2	PRO	4.8
10	1I	47	PHE	4.8
9	82	11	LYS	4.8
14	5A	15	LYS	4.8
4	32	69	GLY	4.8
38	55	8	ARG	4.8
53	O8	50	ARG	4.8
10	1A	66	ARG	4.8
16	7I	23	ASP	4.8
9	8E	121	ARG	4.8
24	3L	16	U	4.8
34	58	74	ARG	4.8
9	82	122	ALA	4.8
31	49	92	VAL	4.8
48	F5	95	LEU	4.7
14	5A	7	ILE	4.7
29	29	54	GLN	4.7
24	1L	74	C	4.7
34	15	109	LYS	4.7
16	7I	10	GLY	4.7
33	69	37	VAL	4.7
31	49	39	ILE	4.7
10	1A	49	VAL	4.7
33	69	11	ASN	4.7
14	5I	59	ALA	4.7
1	1G	1224	G	4.7
7	6E	84	ASN	4.7
24	1L	1	G	4.7
8	72	90	GLY	4.7
33	69	12	LEU	4.7
1	1G	975	A	4.7
32	59	93	GLY	4.7
32	59	170	ARG	4.7
37	45	88	GLY	4.6
16	7I	21	VAL	4.6
45	C5	44	ILE	4.6
9	8E	124	GLN	4.6
14	5A	57	ARG	4.6
16	7I	33	ILE	4.6

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Mol	Chain	Res	Type	RSRZ
34	15	79	PRO	4.6
37	45	32	TYR	4.6
15	6I	62	GLN	4.6
16	7I	2	VAL	4.6
46	D5	172	ALA	4.6
8	7E	4	ASP	4.6
9	8E	118	LYS	4.6
29	21	204	ALA	4.6
14	5A	56	VAL	4.6
9	82	107	ARG	4.6
47	E5	39	ARG	4.6
53	O8	19	ARG	4.6
9	82	124	GLN	4.6
9	8E	120	ARG	4.5
17	8A	32	TYR	4.5
42	95	80	GLN	4.5
46	D5	168	GLU	4.5
20	BA	71	THR	4.5
34	15	73	THR	4.5
55	Q8	27	THR	4.5
20	BA	83	ARG	4.5
39	65	13	ARG	4.5
53	O8	22	ALA	4.5
51	I5	10	VAL	4.5
20	BI	68	LYS	4.5
44	B5	68	ARG	4.5
9	8E	119	ALA	4.5
37	45	99	PRO	4.5
9	8E	109	VAL	4.5
16	7I	20	VAL	4.5
45	C5	29	GLU	4.5
29	29	76	ARG	4.5
46	D5	116	VAL	4.5
36	35	35	HIS	4.5
14	5A	16	PHE	4.5
37	45	41	TRP	4.4
13	4A	26	GLY	4.4
21	1F	24	ARG	4.4
17	8I	35	VAL	4.4
37	45	22	LYS	4.4
38	98	9	LYS	4.4
9	8E	66	ARG	4.4

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Mol	Chain	Res	Type	RSRZ
40	75	99	LEU	4.4
12	3A	129	ALA	4.4
31	49	36	LYS	4.4
46	D5	151	HIS	4.4
53	O8	15	GLU	4.4
37	45	33	GLY	4.4
37	45	64	ILE	4.4
10	1A	50	ILE	4.4
14	5I	30	ALA	4.4
32	59	106	THR	4.4
21	1B	3	LYS	4.4
51	I5	5	ILE	4.4
16	7I	34	GLU	4.4
47	I8	42	GLY	4.4
16	7I	35	LYS	4.3
26	14	2145	C	4.3
9	82	30	GLY	4.3
7	6E	78	ARG	4.3
34	58	109	LYS	4.3
37	45	93	TYR	4.3
32	59	114	VAL	4.3
14	5I	28	GLY	4.3
31	49	139	LEU	4.3
38	55	69	ASP	4.3
21	1B	17	THR	4.3
10	1I	45	ARG	4.3
19	AI	71	LEU	4.3
47	E5	22	GLY	4.3
10	1I	58	ASP	4.3
31	41	26	GLN	4.3
46	D5	79	ARG	4.3
42	95	76	LYS	4.3
8	72	92	ARG	4.3
13	4A	7	VAL	4.3
8	7E	5	PRO	4.3
29	29	141	ILE	4.3
37	45	60	ARG	4.3
31	49	74	LYS	4.3
32	59	105	LEU	4.3
53	O8	21	TYR	4.2
4	3E	209	ARG	4.2
42	95	81	TYR	4.2

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Mol	Chain	Res	Type	RSRZ
14	5I	29	ARG	4.2
53	O8	10	LEU	4.2
9	82	68	GLY	4.2
16	7A	33	ILE	4.2
47	E5	44	ARG	4.2
9	82	75	ASP	4.2
28	19	38	LYS	4.2
9	82	38	GLN	4.2
10	1A	70	ARG	4.2
21	1F	13	ILE	4.2
29	29	77	ILE	4.2
34	15	74	ARG	4.2
32	59	87	LEU	4.2
39	65	7	TYR	4.2
47	E5	42	GLY	4.2
1	1G	1117	G	4.2
10	1I	48	THR	4.2
10	1A	61	GLU	4.2
14	5I	7	ILE	4.2
4	3E	3	ARG	4.2
16	7A	1	MET	4.2
32	59	83	TYR	4.2
51	M8	64	GLY	4.2
1	1G	1250	A	4.1
20	BA	15	ARG	4.1
26	14	2118	U	4.1
39	65	37	ALA	4.1
16	7I	14	ASN	4.1
50	H5	28	LEU	4.1
10	1I	49	VAL	4.1
13	4I	100	GLY	4.1
17	8A	7	THR	4.1
31	41	25	TYR	4.1
32	59	131	VAL	4.1
34	15	81	GLY	4.1
8	72	86	ILE	4.1
20	BA	72	LEU	4.1
31	49	160	VAL	4.1
33	69	21	VAL	4.1
16	7I	36	ILE	4.1
7	62	86	GLN	4.1
40	B8	94	ALA	4.1

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Mol	Chain	Res	Type	RSRZ
1	1G	974	A	4.1
10	1A	34	VAL	4.1
29	29	115	GLY	4.1
29	29	116	VAL	4.1
37	45	105	GLU	4.1
16	7I	26	ARG	4.1
20	BA	18	GLN	4.1
55	Q8	35	GLN	4.1
8	72	88	LYS	4.1
36	35	16	ARG	4.1
38	55	68	ARG	4.1
47	E5	41	ARG	4.1
42	95	75	PHE	4.1
29	29	149	ARG	4.1
41	C8	28	ARG	4.1
47	E5	24	LYS	4.1
26	14	1	G	4.1
14	5A	18	VAL	4.1
20	BA	68	LYS	4.1
20	BA	80	ARG	4.1
29	29	155	LYS	4.1
1	1G	1115	C	4.1
20	BA	14	LYS	4.1
8	72	58	TYR	4.1
26	14	3	U	4.0
39	65	17	ARG	4.0
19	AA	76	PRO	4.0
42	95	77	ALA	4.0
20	BA	75	ASN	4.0
37	45	69	PHE	4.0
9	82	27	THR	4.0
10	1I	6	ILE	4.0
55	Q8	3	LYS	4.0
16	7I	64	ALA	4.0
10	1I	71	LEU	4.0
51	I5	31	ILE	4.0
10	1I	63	PHE	4.0
3	22	177	THR	4.0
26	14	2144	U	4.0
37	45	6	ARG	4.0
10	1I	43	ARG	4.0
20	BI	71	THR	4.0

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Mol	Chain	Res	Type	RSRZ
48	F5	28	GLY	4.0
33	69	6	LEU	4.0
29	29	113	PHE	4.0
36	35	71	VAL	4.0
39	65	14	VAL	4.0
29	29	159	HIS	4.0
8	7E	89	PRO	4.0
10	1A	63	PHE	4.0
16	7I	3	LYS	4.0
16	7I	25	ARG	4.0
28	19	5	LYS	4.0
45	C5	34	LYS	4.0
18	9A	88	LYS	3.9
26	14	2899	G	3.9
9	82	44	VAL	3.9
26	14	2173	A	3.9
47	E5	74	ARG	3.9
31	41	2	PRO	3.9
21	1F	10	ARG	3.9
47	E5	72	ARG	3.9
8	72	22	GLU	3.9
10	1A	44	VAL	3.9
10	1A	68	HIS	3.9
7	62	85	TYR	3.9
8	7E	91	ARG	3.9
24	3L	20	U	3.9
13	4A	65	LYS	3.9
24	1L	3	C	3.9
1	1G	1225	A	3.9
5	42	126	ARG	3.9
46	D5	120	ILE	3.9
8	72	59	LEU	3.9
13	4A	98	VAL	3.9
19	AA	35	SER	3.9
39	65	9	ARG	3.9
20	BI	66	ALA	3.9
32	59	129	THR	3.9
20	BA	76	ALA	3.9
50	H5	17	LYS	3.9
51	I5	32	TYR	3.8
12	3A	69	TYR	3.8
10	1A	43	ARG	3.8

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Mol	Chain	Res	Type	RSRZ
40	B8	99	LEU	3.8
46	H8	113	ALA	3.8
26	1H	2117	A	3.8
55	M5	12	LYS	3.8
39	65	10	ARG	3.8
37	45	12	GLN	3.8
37	45	106	VAL	3.8
5	42	130	ASN	3.8
40	75	104	ASN	3.8
55	M5	44	LYS	3.8
14	5A	53	LEU	3.8
9	82	63	ILE	3.8
26	14	1093	G	3.8
29	21	75	VAL	3.8
55	Q8	26	LYS	3.8
19	AA	84	GLY	3.8
37	45	130	LYS	3.8
16	7I	8	ARG	3.8
41	85	40	PHE	3.8
7	6E	16	LEU	3.8
32	59	168	PRO	3.8
53	O8	12	GLU	3.8
55	Q8	57	ARG	3.8
4	3E	110	PHE	3.8
17	8I	37	LYS	3.7
13	4A	2	ALA	3.7
51	I5	54	GLY	3.7
7	62	16	LEU	3.7
46	D5	173	ALA	3.7
53	K5	43	CYS	3.7
20	BA	29	LYS	3.7
32	51	172	LYS	3.7
41	C8	56	ASP	3.7
53	O8	43	CYS	3.7
24	3L	18	G	3.7
34	15	48	MET	3.7
30	39	97	TYR	3.7
15	6I	60	VAL	3.7
47	E5	69	PHE	3.7
16	7A	7	ALA	3.7
9	82	40	LEU	3.7
47	E5	77	ARG	3.7

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Mol	Chain	Res	Type	RSRZ
24	1L	70	G	3.7
34	58	84	LYS	3.7
1	13	134	A	3.7
16	7A	59	TRP	3.7
17	8A	24	GLU	3.7
8	7E	92	ARG	3.7
19	AA	53	ASN	3.7
12	3A	5	PRO	3.7
15	6A	31	LEU	3.7
29	29	136	ARG	3.7
28	19	212	SER	3.7
34	15	116	LEU	3.7
48	F5	3	LYS	3.7
51	I5	28	LYS	3.7
17	8A	71	PHE	3.7
44	B5	69	TYR	3.7
34	15	77	GLY	3.6
3	22	178	LEU	3.6
36	35	62	LEU	3.6
35	25	33	ALA	3.6
26	14	2111	C	3.6
9	8E	128	ARG	3.6
1	1G	1287	A	3.6
36	35	45	LEU	3.6
39	65	20	ARG	3.6
39	65	112	PHE	3.6
36	35	149	GLU	3.6
20	BI	79	ARG	3.6
44	B5	33	LYS	3.6
45	C5	48	ALA	3.6
10	1I	61	GLU	3.6
29	29	160	TYR	3.6
8	72	4	ASP	3.6
9	8E	75	ASP	3.6
9	82	32	ASP	3.6
9	82	82	ALA	3.6
10	1I	50	ILE	3.6
36	35	107	LYS	3.6
55	M5	59	LYS	3.6
13	4A	73	GLU	3.6
26	14	1095	A	3.6
34	15	76	SER	3.6

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Mol	Chain	Res	Type	RSRZ
51	I5	24	THR	3.6
15	6I	48	LYS	3.6
29	29	59	VAL	3.6
55	Q8	21	LYS	3.6
7	62	2	ALA	3.6
15	6I	65	ARG	3.6
4	3E	2	GLY	3.6
31	49	12	TYR	3.6
5	4E	123	LEU	3.6
12	3I	20	LYS	3.6
28	19	206	LEU	3.6
7	62	13	GLN	3.6
1	1G	825	G	3.6
53	K5	35	GLU	3.6
46	D5	163	LEU	3.6
51	I5	40	HIS	3.6
1	1G	1249	C	3.6
9	82	102	LEU	3.6
31	49	94	LEU	3.6
29	21	76	ARG	3.6
8	72	93	VAL	3.6
14	5I	2	ALA	3.6
53	K5	30	THR	3.6
55	M5	3	LYS	3.6
10	1A	45	ARG	3.6
47	E5	19	LYS	3.6
17	8I	32	TYR	3.6
9	82	9	ARG	3.5
51	M8	59	PHE	3.5
8	72	136	GLU	3.5
12	3I	91	LYS	3.5
2	1E	187	LEU	3.5
14	5A	12	ARG	3.5
20	BA	63	ILE	3.5
20	BA	13	LEU	3.5
21	1B	4	GLY	3.5
26	14	2170	A	3.5
37	45	8	LYS	3.5
1	13	230	G	3.5
9	82	35	GLU	3.5
31	49	157	ILE	3.5
4	32	209	ARG	3.5

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Mol	Chain	Res	Type	RSRZ
7	62	32	ARG	3.5
9	82	72	GLY	3.5
40	B8	104	ASN	3.5
38	55	13	HIS	3.5
16	7A	26	ARG	3.5
20	BI	29	LYS	3.5
20	BA	84	LEU	3.5
9	8E	125	TYR	3.5
4	32	83	SER	3.5
9	8E	7	THR	3.5
47	E5	55	ARG	3.5
4	3E	96	LEU	3.5
15	6I	28	GLN	3.5
36	35	64	LYS	3.5
40	75	6	LEU	3.5
9	82	62	TYR	3.5
53	K5	44	ARG	3.5
53	O8	47	THR	3.5
21	1F	9	ARG	3.5
2	12	197	VAL	3.5
47	I8	46	LYS	3.5
39	65	32	LEU	3.5
48	F5	10	LYS	3.5
16	7I	13	HIS	3.5
48	J8	95	LEU	3.5
37	45	7	MET	3.5
2	1E	188	ALA	3.5
10	1I	42	THR	3.5
37	45	14	ARG	3.5
46	D5	176	PRO	3.5
17	8A	4	LYS	3.5
34	15	83	LYS	3.5
13	4A	104	ARG	3.5
20	BA	25	ARG	3.5
41	C8	26	GLY	3.5
47	I8	40	GLN	3.5
34	58	71	ILE	3.5
17	8A	68	ARG	3.5
9	8E	112	LYS	3.5
4	3E	137	SER	3.5
10	1I	59	SER	3.5
32	59	159	GLU	3.5

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Mol	Chain	Res	Type	RSRZ
4	3E	139	ARG	3.5
8	72	84	ARG	3.5
39	65	35	ILE	3.5
47	E5	73	GLY	3.5
33	61	117	GLU	3.4
29	21	90	THR	3.4
52	J5	11	THR	3.4
53	O8	52	VAL	3.4
20	BA	23	ARG	3.4
33	61	113	ARG	3.4
17	8A	59	ILE	3.4
48	F5	94	LEU	3.4
3	22	206	GLU	3.4
8	7E	88	LYS	3.4
16	7I	68	ASP	3.4
34	15	72	TYR	3.4
20	BA	26	ASN	3.4
13	4I	104	ARG	3.4
2	12	118	LEU	3.4
16	7A	32	TYR	3.4
4	32	70	ILE	3.4
5	42	45	PHE	3.4
1	1G	1028(B)	C	3.4
14	5A	6	LEU	3.4
12	3A	48	PRO	3.4
16	7A	31	LYS	3.4
41	C8	29	SER	3.4
46	D5	9	TYR	3.4
9	82	33	PHE	3.4
37	45	11	LYS	3.4
26	1H	529	A	3.4
45	C5	5	MET	3.4
42	95	91	TYR	3.4
19	AA	83	HIS	3.4
20	BI	8	ARG	3.4
29	29	162	ALA	3.4
14	5A	47	LEU	3.4
15	6A	57	LEU	3.4
20	BA	65	LYS	3.4
39	65	6	ALA	3.4
51	M8	56	VAL	3.4
9	82	105	ASP	3.4

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Mol	Chain	Res	Type	RSRZ
1	1G	1202	G	3.4
37	45	35	VAL	3.4
14	5A	13	THR	3.4
37	45	92	GLY	3.4
34	15	1	MET	3.4
20	BA	66	ALA	3.4
45	C5	60	PHE	3.4
9	8E	42	ARG	3.4
48	F5	11	ARG	3.4
2	12	115	LEU	3.3
16	7I	37	GLY	3.3
38	55	10	LEU	3.3
1	1G	1251	A	3.3
4	3E	168	ARG	3.3
5	4E	122	GLU	3.3
28	19	217	ARG	3.3
2	12	152	PHE	3.3
41	85	52	ARG	3.3
50	H5	10	LYS	3.3
9	82	73	GLN	3.3
26	14	2319	G	3.3
43	A5	92	ARG	3.3
47	I8	57	PHE	3.3
13	4I	97	PRO	3.3
7	62	4	ARG	3.3
26	14	2799	A	3.3
36	78	150	ALA	3.3
35	25	32	TYR	3.3
38	98	69	ASP	3.3
9	8E	10	ARG	3.3
9	82	43	ALA	3.3
9	8E	123	PRO	3.3
50	H5	12	PRO	3.3
31	49	37	VAL	3.3
32	59	37	VAL	3.3
55	M5	4	MET	3.3
26	14	1537	C	3.3
29	29	154	LYS	3.3
24	3L	19	G	3.3
40	75	50	ILE	3.3
26	14	1026	U	3.3
41	C8	27	LEU	3.3

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Mol	Chain	Res	Type	RSRZ
29	29	122	PHE	3.3
46	D5	118	GLN	3.3
1	1G	1223	C	3.3
22	1K	75	C	3.3
31	49	28	VAL	3.3
37	45	97	VAL	3.3
10	1I	98	ILE	3.3
29	29	157	ALA	3.3
7	62	36	LYS	3.3
19	AA	12	ASP	3.3
20	BA	104	LEU	3.3
34	15	78	TYR	3.3
31	41	23	PHE	3.3
45	C5	24	VAL	3.3
31	49	155	MET	3.2
38	98	7	GLY	3.2
1	1G	1001	G	3.2
26	14	2148	G	3.2
52	J5	2	ALA	3.2
37	88	104	PHE	3.2
9	8E	78	LYS	3.2
9	82	78	LYS	3.2
19	AA	36	ARG	3.2
20	BI	23	ARG	3.2
34	15	46	VAL	3.2
13	4A	92	HIS	3.2
37	45	73	PRO	3.2
14	5A	44	LEU	3.2
39	65	4	LEU	3.2
20	BA	79	ARG	3.2
29	29	75	VAL	3.2
16	7I	16	HIS	3.2
19	AA	69	HIS	3.2
29	29	163	GLU	3.2
9	8E	102	LEU	3.2
9	8E	37	PHE	3.2
37	45	29	PHE	3.2
5	4E	11	ILE	3.2
5	4E	89	ILE	3.2
16	7I	15	PRO	3.2
43	A5	82	LEU	3.2
31	49	159	VAL	3.2

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Mol	Chain	Res	Type	RSRZ
41	85	47	TYR	3.2
42	D8	81	TYR	3.2
15	6A	59	MET	3.2
17	8A	8	GLY	3.2
42	95	83	ARG	3.2
39	65	11	LYS	3.2
5	42	24	ARG	3.2
29	21	79	ARG	3.2
38	98	67	LEU	3.2
46	D5	147	GLY	3.2
12	3A	16	GLU	3.2
7	6E	32	ARG	3.2
12	3A	15	ARG	3.2
31	49	96	ARG	3.2
32	59	97	ARG	3.2
8	72	17	THR	3.2
30	39	82	ILE	3.2
19	AA	70	LYS	3.2
20	BA	74	LYS	3.2
32	59	153	LYS	3.2
45	C5	66	PRO	3.2
4	3E	135	LEU	3.2
50	H5	19	GLN	3.2
51	I5	46	GLN	3.2
21	1F	23	PRO	3.2
1	13	1032(A)	G	3.2
47	E5	45	PHE	3.2
13	4A	5	ALA	3.2
47	I8	41	ARG	3.2
1	1G	1116	C	3.2
5	4E	81	GLU	3.2
16	7I	38	TYR	3.2
50	H5	15	TYR	3.2
15	6I	31	LEU	3.2
29	29	112	GLY	3.2
31	49	175	LEU	3.2
20	BA	78	ALA	3.2
13	4A	87	TYR	3.2
36	35	47	ASP	3.2
37	45	129	THR	3.2
41	85	6	THR	3.2
10	1I	57	LYS	3.1

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Mol	Chain	Res	Type	RSRZ
13	4I	96	LEU	3.1
33	6I	118	LYS	3.1
3	22	10	PHE	3.1
9	8E	9	ARG	3.1
34	15	115	ARG	3.1
48	F5	93	GLU	3.1
20	BI	20	LEU	3.1
46	D5	76	LEU	3.1
41	85	3	ARG	3.1
9	8E	14	VAL	3.1
33	69	18	VAL	3.1
10	1I	40	LEU	3.1
29	29	105	THR	3.1
32	51	171	LEU	3.1
53	O8	37	ARG	3.1
32	59	161	GLY	3.1
42	95	70	ILE	3.1
55	Q8	48	PHE	3.1
37	45	80	GLU	3.1
48	J8	70	VAL	3.1
9	82	31	GLN	3.1
26	1H	2116	G	3.1
30	31	40	GLN	3.1
11	2A	126	ARG	3.1
28	19	211	ARG	3.1
29	29	111	ARG	3.1
8	7E	131	GLY	3.1
8	72	55	GLY	3.1
10	1A	11	PHE	3.1
14	5A	54	PRO	3.1
21	1F	20	LYS	3.1
53	O8	33	LYS	3.1
11	2I	12	ARG	3.1
26	14	2114	A	3.1
19	AA	13	ASP	3.1
32	59	112	PRO	3.1
48	F5	32	LYS	3.1
53	O8	48	VAL	3.1
51	I5	9	LEU	3.1
2	1E	27	LYS	3.1
15	6I	72	ARG	3.1
34	15	86	PRO	3.1

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Mol	Chain	Res	Type	RSRZ
5	42	88	LYS	3.1
38	55	70	LEU	3.1
26	14	2125	G	3.1
34	58	115	ARG	3.1
55	M5	13	ARG	3.1
47	E5	23	VAL	3.1
47	E5	38	VAL	3.1
47	E5	53	MET	3.1
7	62	88	PRO	3.1
10	1A	41	PRO	3.1
4	32	19	LEU	3.1
8	72	112	LEU	3.1
16	7I	67	THR	3.1
37	45	98	LYS	3.1
15	6I	27	VAL	3.1
55	Q8	5	LYS	3.1
46	D5	125	LEU	3.1
20	BA	17	ARG	3.1
20	BA	67	ALA	3.1
28	19	37	LEU	3.1
37	45	140	ALA	3.1
9	82	77	ILE	3.1
9	82	28	VAL	3.1
1	13	1367	C	3.1
7	6E	5	ARG	3.1
7	62	38	LEU	3.1
39	65	87	PHE	3.0
51	I5	11	PRO	3.0
8	7E	93	VAL	3.0
8	72	91	ARG	3.0
13	4A	88	ARG	3.0
34	15	44	PRO	3.0
16	7A	64	ALA	3.0
1	1G	876	G	3.0
19	AA	40	ILE	3.0
55	M5	41	ILE	3.0
47	I8	53	MET	3.0
28	19	235	GLY	3.0
39	65	18	ILE	3.0
47	I8	52	GLY	3.0
14	5I	31	ARG	3.0
7	6E	35	LYS	3.0

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Mol	Chain	Res	Type	RSRZ
32	59	85	LYS	3.0
38	55	14	SER	3.0
48	F5	92	LYS	3.0
1	1G	1357	A	3.0
4	3E	93	PHE	3.0
17	8I	27	PHE	3.0
55	Q8	6	THR	3.0
29	21	4	ILE	3.0
38	55	2	ARG	3.0
7	62	42	ILE	3.0
21	1F	4	GLY	3.0
26	14	1046	A	3.0
4	3E	21	LEU	3.0
29	21	195	LEU	3.0
10	1I	54	PHE	3.0
15	6A	68	ARG	3.0
10	1A	10	GLY	3.0
12	3A	68	ALA	3.0
34	58	73	THR	3.0
20	BA	62	LEU	3.0
14	5I	32	SER	3.0
19	AA	49	ILE	3.0
29	29	192	ASN	3.0
41	C8	34	LYS	3.0
31	49	161	THR	3.0
41	85	20	LEU	3.0
55	Q8	62	LEU	3.0
26	1H	654(K)	C	3.0
32	51	3	ARG	3.0
37	45	39	PRO	3.0
1	1G	1248	A	3.0
10	1I	55	LYS	3.0
39	65	93	LYS	3.0
8	72	18	ARG	3.0
20	BI	14	LYS	3.0
47	E5	40	GLN	3.0
55	M5	26	LYS	3.0
26	14	2138	C	3.0
28	19	55	GLY	3.0
1	1G	1236	A	3.0
1	1G	1285	A	3.0
2	1E	31	TYR	3.0

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Mol	Chain	Res	Type	RSRZ
4	32	122	ARG	3.0
9	82	79	LEU	3.0
10	1A	51	ARG	3.0
13	4A	94	ARG	3.0
20	BI	80	ARG	3.0
34	15	82	LEU	3.0
36	35	42	SER	3.0
39	65	57	LYS	3.0
50	H5	20	LYS	3.0
1	1G	1358	U	3.0
41	C8	25	TRP	3.0
1	1G	1367	C	3.0
16	7A	6	LEU	3.0
29	29	152	LYS	3.0
10	1I	67	THR	3.0
30	39	92	PRO	3.0
5	42	14	ARG	2.9
48	F5	21	ARG	2.9
29	21	72	VAL	2.9
45	C5	49	VAL	2.9
8	72	56	LYS	2.9
14	5I	17	LYS	2.9
45	C5	63	LYS	2.9
1	13	1354	C	2.9
36	35	13	ASN	2.9
20	BA	73	HIS	2.9
4	3E	138	TYR	2.9
46	D5	51	ALA	2.9
33	69	16	GLY	2.9
33	69	34	GLY	2.9
36	35	46	LYS	2.9
20	BI	67	ALA	2.9
26	14	1536	A	2.9
32	59	141	VAL	2.9
55	M5	34	TRP	2.9
8	72	25	ASP	2.9
49	K8	69	ARG	2.9
14	5A	42	ILE	2.9
29	29	137	HIS	2.9
37	45	20	ALA	2.9
37	45	37	LEU	2.9
42	D8	38	LEU	2.9

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Mol	Chain	Res	Type	RSRZ
15	6I	54	ARG	2.9
47	I8	45	PHE	2.9
8	72	6	ILE	2.9
2	12	28	PHE	2.9
8	7E	85	ARG	2.9
26	14	2318	G	2.9
37	45	5	ARG	2.9
40	75	98	LYS	2.9
45	C5	64	GLU	2.9
37	45	100	GLY	2.9
5	42	119	LEU	2.9
14	5I	57	ARG	2.9
37	45	120	ILE	2.9
5	42	84	PHE	2.9
12	3A	47	LYS	2.9
28	19	39	LYS	2.9
29	29	203	LYS	2.9
36	35	51	PHE	2.9
8	7E	90	GLY	2.9
17	8A	11	VAL	2.9
20	BI	17	ARG	2.9
1	13	1325	C	2.9
26	14	1661	G	2.9
51	I5	42	PHE	2.9
13	4A	105	THR	2.9
55	M5	23	VAL	2.9
2	12	40	HIS	2.9
20	BI	16	HIS	2.9
32	59	130	ARG	2.9
55	M5	2	PRO	2.9
1	13	1032	A	2.9
8	7E	133	LEU	2.9
20	BI	63	ILE	2.9
36	35	38	GLN	2.9
9	8E	105	ASP	2.9
33	69	20	ASP	2.9
15	6I	63	ARG	2.9
55	M5	8	LYS	2.9
1	13	311	C	2.9
8	72	61	VAL	2.9
12	3I	90	VAL	2.9
28	19	257	LEU	2.9

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Mol	Chain	Res	Type	RSRZ
31	49	27	ASN	2.9
1	1G	977	A	2.9
11	2A	25	TYR	2.9
31	49	95	ARG	2.9
29	29	145	LYS	2.9
51	I5	7	PRO	2.9
1	13	229	U	2.8
1	1G	972	C	2.8
10	1A	42	THR	2.8
29	29	107	THR	2.8
30	39	72	ARG	2.8
55	M5	15	LYS	2.8
55	M5	21	LYS	2.8
1	13	111	G	2.8
12	3A	85	ILE	2.8
29	29	156	MET	2.8
12	3I	17	LYS	2.8
36	35	30	THR	2.8
37	45	40	ALA	2.8
51	I5	30	GLU	2.8
26	14	2164	C	2.8
16	7A	9	PHE	2.8
20	BA	22	ARG	2.8
40	B8	101	PHE	2.8
1	1G	1373	G	2.8
24	3L	60	U	2.8
1	1G	1354	C	2.8
13	4A	25	ILE	2.8
30	39	93	LYS	2.8
39	65	12	PHE	2.8
19	AI	38	SER	2.8
26	14	2713	A	2.8
20	BI	9	ASN	2.8
1	13	112	G	2.8
1	1G	966	G	2.8
29	29	24	THR	2.8
46	D5	78	LYS	2.8
39	65	91	PRO	2.8
20	BI	106	ALA	2.8
19	AI	78	ARG	2.8
28	19	4	LYS	2.8
28	19	155	LEU	2.8

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Mol	Chain	Res	Type	RSRZ
36	78	30	THR	2.8
46	D5	141	VAL	2.8
29	21	147	PRO	2.8
40	75	100	TYR	2.8
42	95	78	LYS	2.8
14	5I	60	SER	2.8
16	7I	11	SER	2.8
1	13	1364	U	2.8
4	3E	111	ALA	2.8
11	2I	126	ARG	2.8
20	BA	81	LYS	2.8
38	98	17	ARG	2.8
2	12	135	GLN	2.8
14	5A	27	CYS	2.8
15	6A	62	GLN	2.8
8	72	12	ARG	2.8
9	82	76	ALA	2.8
7	6E	34	GLY	2.8
13	4A	90	LEU	2.8
38	55	21	TYR	2.8
41	C8	35	ALA	2.8
53	O8	34	LEU	2.8
32	59	162	ILE	2.8
12	3A	8	ASN	2.8
32	59	86	GLU	2.8
12	3A	98	TYR	2.8
28	19	177	LEU	2.8
31	49	152	LEU	2.8
2	12	163	PHE	2.8
9	82	74	ILE	2.8
14	5I	8	GLU	2.8
20	BI	21	LYS	2.8
36	78	64	LYS	2.8
55	Q8	13	ARG	2.8
46	D5	149	SER	2.8
1	1G	1226	C	2.8
8	7E	112	LEU	2.8
28	19	254	THR	2.8
1	13	1033	G	2.8
41	85	56	ASP	2.8
47	I8	76	GLY	2.8
8	72	111	ILE	2.8

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Mol	Chain	Res	Type	RSRZ
16	7I	5	ARG	2.8
8	7E	9	MET	2.8
43	A5	81	ALA	2.8
53	O8	40	CYS	2.8
20	BI	38	LYS	2.8
53	K5	31	PRO	2.8
41	85	25	TRP	2.7
4	32	11	LEU	2.7
1	13	609	A	2.7
14	5I	3	ARG	2.7
29	21	152	LYS	2.7
14	5I	37	PHE	2.7
39	65	58	LEU	2.7
40	B8	93	ARG	2.7
40	B8	114	LEU	2.7
26	14	2113	U	2.7
9	8E	122	ALA	2.7
29	29	51	PHE	2.7
29	29	143	ASN	2.7
8	72	131	GLY	2.7
20	BI	69	GLY	2.7
31	49	89	GLY	2.7
5	4E	129	ILE	2.7
29	21	55	ASN	2.7
41	C8	37	GLU	2.7
14	5I	22	THR	2.7
24	3L	35	A	2.7
19	AI	75	ALA	2.7
1	13	43	C	2.7
26	14	2142	C	2.7
17	8A	37	LYS	2.7
36	35	33	ARG	2.7
26	1H	2476	A	2.7
29	21	67	PHE	2.7
34	15	117	PHE	2.7
32	59	113	VAL	2.7
8	7E	83	ILE	2.7
16	7A	25	ARG	2.7
20	BI	33	ILE	2.7
52	J5	10	LYS	2.7
54	L5	14	LYS	2.7
55	M5	57	ARG	2.7

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Mol	Chain	Res	Type	RSRZ
11	2A	118	GLY	2.7
3	22	30	ARG	2.7
21	1B	12	LYS	2.7
34	15	114	ARG	2.7
26	14	2112	G	2.7
19	AA	38	SER	2.7
47	E5	57	PHE	2.7
7	6E	33	ASP	2.7
30	39	96	ASP	2.7
53	O8	36	LEU	2.7
1	13	815	A	2.7
1	1G	1033	G	2.7
17	8A	38	ARG	2.7
17	8A	69	LYS	2.7
29	29	11	MET	2.7
39	65	19	LYS	2.7
48	F5	5	CYS	2.7
36	35	27	HIS	2.7
55	Q8	28	GLY	2.7
20	BA	12	ALA	2.7
51	I5	44	THR	2.7
5	42	123	LEU	2.7
55	M5	61	LEU	2.7
36	35	68	GLN	2.7
4	3E	124	GLY	2.7
45	C5	23	ARG	2.7
5	42	81	GLU	2.7
9	8E	65	VAL	2.7
31	49	137	GLU	2.7
29	29	121	ASN	2.7
37	45	74	TYR	2.7
37	45	89	ASN	2.7
1	13	754	C	2.7
5	4E	88	LYS	2.7
26	14	2161	C	2.7
12	3A	18	VAL	2.7
28	11	210	GLY	2.7
29	29	167	VAL	2.7
7	6E	85	TYR	2.7
12	3A	7	ILE	2.7
55	M5	58	ILE	2.7
20	BA	21	LYS	2.7

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Mol	Chain	Res	Type	RSRZ
31	41	34	LEU	2.7
38	55	65	LEU	2.7
4	3E	201	GLN	2.7
9	8E	101	PHE	2.7
37	45	30	GLY	2.7
55	Q8	43	GLN	2.7
2	12	136	VAL	2.7
17	8I	28	PRO	2.7
31	49	135	LEU	2.6
9	82	39	GLY	2.6
3	2E	193	TYR	2.6
5	42	13	ILE	2.6
16	7I	42	ARG	2.6
17	8A	36	ILE	2.6
29	29	191	PRO	2.6
46	D5	171	ILE	2.6
47	E5	18	ALA	2.6
1	13	742	G	2.6
29	29	140	SER	2.6
2	12	14	GLY	2.6
45	C5	58	GLY	2.6
7	62	39	ALA	2.6
8	72	13	ILE	2.6
12	3I	5	PRO	2.6
28	19	246	PRO	2.6
1	1G	879	C	2.6
1	13	1031	G	2.6
8	7E	132	GLU	2.6
26	14	1998	G	2.6
35	68	1	MET	2.6
55	M5	35	GLN	2.6
9	8E	40	LEU	2.6
17	8A	60	ILE	2.6
29	21	52	LEU	2.6
29	21	187	ALA	2.6
48	F5	63	ALA	2.6
34	15	37	LYS	2.6
38	98	14	SER	2.6
41	85	29	SER	2.6
20	BA	64	ASP	2.6
8	72	10	LEU	2.6
10	1I	8	LEU	2.6

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Mol	Chain	Res	Type	RSRZ
15	6I	87	ILE	2.6
19	AA	71	LEU	2.6
39	65	36	TYR	2.6
53	O8	11	LEU	2.6
26	14	944	G	2.6
53	K5	38	LYS	2.6
36	35	26	GLY	2.6
38	55	71	GLN	2.6
33	69	5	LEU	2.6
48	F5	90	ILE	2.6
2	12	131	PRO	2.6
20	BI	22	ARG	2.6
29	29	119	ARG	2.6
31	49	23	PHE	2.6
1	1G	1353	G	2.6
46	D5	96	VAL	2.6
55	M5	14	VAL	2.6
34	15	90	MET	2.6
28	19	147	LEU	2.6
4	32	68	TYR	2.6
28	19	247	ALA	2.6
36	35	108	LYS	2.6
41	C8	30	LYS	2.6
15	6A	25	THR	2.6
17	8I	91	ARG	2.6
29	21	91	VAL	2.6
29	29	104	VAL	2.6
39	65	28	VAL	2.6
5	4E	13	ILE	2.6
10	1I	38	ILE	2.6
12	3A	13	LYS	2.6
45	C5	61	ILE	2.6
20	BI	59	ALA	2.6
16	7A	69	THR	2.6
13	4I	98	VAL	2.6
17	8I	29	HIS	2.6
29	29	132	HIS	2.6
37	45	94	VAL	2.6
29	21	154	LYS	2.6
5	42	15	ARG	2.6
46	D5	164	ALA	2.6
31	49	15	VAL	2.6

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Mol	Chain	Res	Type	RSRZ
46	D5	169	GLU	2.6
8	72	5	PRO	2.6
15	6I	66	LEU	2.6
17	8I	31	LEU	2.6
41	85	49	HIS	2.6
28	19	90	ALA	2.6
42	D8	75	PHE	2.6
11	2I	124	LYS	2.6
29	29	133	LYS	2.6
34	15	45	ASN	2.6
41	C8	118	GLY	2.6
40	B8	115	ARG	2.6
46	D5	170	THR	2.6
55	M5	30	ARG	2.6
1	1G	963	G	2.6
20	BA	27	LYS	2.6
26	14	2167	U	2.6
34	15	70	LYS	2.6
48	J8	69	LYS	2.6
55	M5	29	LYS	2.6
10	1I	44	VAL	2.6
16	7I	61	SER	2.6
55	M5	56	GLU	2.6
38	55	17	ARG	2.6
39	A8	9	ARG	2.6
20	BA	16	HIS	2.6
14	5I	10	ALA	2.5
32	59	104	GLU	2.5
34	58	75	TYR	2.5
1	1G	1362	C	2.5
16	7I	53	VAL	2.5
26	14	2	G	2.5
29	21	199	ARG	2.5
29	29	106	GLY	2.5
50	H5	9	VAL	2.5
32	59	57	ASP	2.5
17	8I	34	LYS	2.5
37	45	63	LYS	2.5
36	78	110	TYR	2.5
2	1E	11	LEU	2.5
34	58	76	SER	2.5
38	98	10	LEU	2.5

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Mol	Chain	Res	Type	RSRZ
39	65	73	LEU	2.5
8	72	9	MET	2.5
12	3A	23	LYS	2.5
7	6E	86	GLN	2.5
37	45	53	ALA	2.5
48	F5	62	VAL	2.5
47	E5	59	LEU	2.5
41	C8	40	PHE	2.5
1	1G	994	A	2.5
30	39	95	ARG	2.5
41	85	2	PRO	2.5
55	Q8	7	HIS	2.5
55	Q8	61	LEU	2.5
37	45	101	ARG	2.5
55	M5	46	ARG	2.5
9	82	17	VAL	2.5
14	5I	58	LYS	2.5
16	7A	63	GLY	2.5
21	1F	11	GLY	2.5
26	14	960	A	2.5
32	59	160	LYS	2.5
35	25	65	THR	2.5
37	45	96	VAL	2.5
48	F5	69	LYS	2.5
55	M5	27	THR	2.5
47	I8	39	ARG	2.5
21	1F	5	ASP	2.5
43	A5	94	ASP	2.5
8	7E	95	VAL	2.5
36	78	107	LYS	2.5
35	68	122	LEU	2.5
37	45	125	LEU	2.5
47	I8	59	LEU	2.5
50	H5	8	LEU	2.5
5	42	25	ARG	2.5
14	5I	13	THR	2.5
9	82	18	PHE	2.5
34	15	51	PHE	2.5
40	75	101	PHE	2.5
32	59	41	MET	2.5
31	49	13	GLU	2.5
38	98	15	SER	2.5

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Mol	Chain	Res	Type	RSRZ
38	55	7	GLY	2.5
55	Q8	2	PRO	2.5
1	1G	1029	G	2.5
16	7I	60	LEU	2.5
48	F5	66	HIS	2.5
47	E5	12	ASN	2.5
1	1G	983	A	2.5
26	14	571	A	2.5
4	3E	203	VAL	2.5
4	32	76	ARG	2.5
13	4A	96	LEU	2.5
15	6A	65	ARG	2.5
28	19	54	ARG	2.5
28	19	182	LEU	2.5
36	35	28	GLY	2.5
36	35	50	ARG	2.5
36	35	53	GLY	2.5
48	F5	64	ALA	2.5
8	72	31	PHE	2.5
29	29	120	TRP	2.5
33	69	40	THR	2.5
34	15	118	LYS	2.5
37	45	67	ARG	2.5
44	B5	54	VAL	2.5
1	13	389	A	2.5
36	35	48	PRO	2.5
4	3E	115	ARG	2.5
36	35	32	THR	2.5
51	I5	55	ARG	2.5
3	22	155	GLY	2.5
11	2I	25	TYR	2.5
12	3A	84	LEU	2.5
16	7I	49	LEU	2.5
28	19	2	ALA	2.5
28	19	18	VAL	2.5
43	E8	23	LEU	2.5
1	13	1366	C	2.5
34	58	85	ILE	2.5
43	A5	103	ILE	2.5
12	3I	11	VAL	2.5
31	41	27	ASN	2.5
17	8I	98	LEU	2.5

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Mol	Chain	Res	Type	RSRZ
37	45	102	VAL	2.5
41	C8	38	THR	2.5
45	C5	39	VAL	2.5
36	35	31	ALA	2.5
41	85	30	LYS	2.5
10	1A	6	ILE	2.5
48	F5	8	SER	2.5
52	N8	55	ARG	2.5
1	1G	132	C	2.4
9	8E	113	LYS	2.4
43	E8	113	LYS	2.4
32	59	157	TYR	2.4
53	K5	28	ARG	2.4
29	29	127	ASP	2.4
9	8E	70	LYS	2.4
15	6A	5	LYS	2.4
1	1G	1032(B)	G	2.4
1	1G	1369	C	2.4
24	3K	65	G	2.4
40	75	5	ALA	2.4
31	49	93	THR	2.4
13	4A	111	LYS	2.4
29	29	129	HIS	2.4
8	72	87	SER	2.4
48	F5	70	VAL	2.4
31	49	178	PHE	2.4
39	65	92	TYR	2.4
8	7E	6	ILE	2.4
8	7E	136	GLU	2.4
1	1G	814	A	2.4
1	1G	969	A	2.4
12	3A	46	LYS	2.4
7	6E	80	VAL	2.4
8	7E	84	ARG	2.4
29	29	125	GLY	2.4
2	1E	152	PHE	2.4
4	32	79	PHE	2.4
28	11	247	ALA	2.4
55	M5	24	ALA	2.4
36	35	29	LYS	2.4
32	59	128	PRO	2.4
32	59	35	VAL	2.4

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Mol	Chain	Res	Type	RSRZ
1	1G	964	A	2.4
1	1G	1363	A	2.4
26	14	2116	G	2.4
7	62	103	TRP	2.4
9	82	46	ALA	2.4
4	3E	5	ILE	2.4
48	F5	42	GLN	2.4
15	6I	67	LEU	2.4
17	8A	6	LEU	2.4
10	1I	10	GLY	2.4
48	F5	36	GLY	2.4
5	4E	45	PHE	2.4
26	14	4	C	2.4
46	D5	148	ASP	2.4
1	13	262	A	2.4
29	29	204	ALA	2.4
5	4E	101	ILE	2.4
29	29	108	SER	2.4
43	E8	24	ILE	2.4
31	41	75	LYS	2.4
35	68	66	LYS	2.4
40	B8	92	GLY	2.4
8	72	135	CYS	2.4
46	D5	5	LEU	2.4
47	E5	70	GLN	2.4
48	F5	91	LYS	2.4
51	I5	47	GLN	2.4
1	1G	1220	G	2.4
17	8I	71	PHE	2.4
37	45	4	PRO	2.4
9	82	16	ARG	2.4
28	19	183	ARG	2.4
37	45	25	ASP	2.4
48	F5	20	ARG	2.4
1	13	307	C	2.4
26	14	1660	C	2.4
34	58	46	VAL	2.4
34	58	82	LEU	2.4
47	I8	37	LEU	2.4
1	1G	1289	A	2.4
32	59	109	PHE	2.4
41	C8	32	PHE	2.4

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Mol	Chain	Res	Type	RSRZ
12	3A	94	PRO	2.4
1	13	1001	G	2.4
37	45	71	ASP	2.4
31	49	3	LEU	2.4
29	21	198	VAL	2.4
30	39	1	MET	2.4
32	59	84	SER	2.4
37	45	15	GLY	2.4
48	J8	60	PHE	2.4
1	13	135	C	2.4
15	6I	68	ARG	2.4
4	32	77	ASN	2.4
55	M5	54	GLU	2.4
55	M5	47	LYS	2.4
28	11	111	LEU	2.4
29	21	167	VAL	2.4
1	13	1343	G	2.4
37	45	13	GLN	2.4
47	E5	52	GLY	2.4
8	72	85	ARG	2.4
14	5I	12	ARG	2.4
5	4E	128	PRO	2.4
12	3A	71	PRO	2.4
26	14	2691	C	2.4
37	45	121	ALA	2.4
4	32	108	LEU	2.4
17	8A	22	LEU	2.4
41	85	18	LEU	2.4
1	1G	262	A	2.3
16	7I	48	TRP	2.3
17	8A	9	VAL	2.3
32	59	125	VAL	2.3
15	6I	58	MET	2.3
11	2A	123	LYS	2.3
42	95	84	LYS	2.3
10	1A	38	ILE	2.3
29	29	134	ILE	2.3
31	49	19	LEU	2.3
48	J8	90	ILE	2.3
1	13	1362	C	2.3
9	8E	41	VAL	2.3
15	6I	50	HIS	2.3

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Mol	Chain	Res	Type	RSRZ
19	AA	77	THR	2.3
40	B8	112	ARG	2.3
37	45	19	GLY	2.3
42	95	93	GLU	2.3
10	1A	37	PRO	2.3
12	3A	100	ILE	2.3
33	69	39	ALA	2.3
36	35	105	LEU	2.3
55	M5	50	LEU	2.3
9	82	41	VAL	2.3
26	1H	615	G	2.3
37	45	128	LYS	2.3
39	65	60	GLY	2.3
1	1G	984	C	2.3
1	1G	1320	C	2.3
16	7A	34	GLU	2.3
12	3A	93	LEU	2.3
15	6I	34	LEU	2.3
26	14	764	A	2.3
34	58	114	ARG	2.3
49	K8	66	GLU	2.3
53	O8	35	GLU	2.3
4	3E	102	ASP	2.3
1	13	309	G	2.3
1	1G	824	C	2.3
1	1G	1321	C	2.3
15	6I	64	ARG	2.3
39	A8	3	ARG	2.3
5	4E	131	ILE	2.3
20	BI	76	ALA	2.3
29	29	78	LEU	2.3
32	59	145	ALA	2.3
36	78	45	LEU	2.3
37	45	44	ALA	2.3
41	C8	22	LYS	2.3
55	Q8	51	ALA	2.3
12	3I	45	PRO	2.3
16	7A	2	VAL	2.3
17	8A	23	VAL	2.3
32	59	49	VAL	2.3
1	13	1349	A	2.3
29	29	142	GLY	2.3

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Mol	Chain	Res	Type	RSRZ
5	4E	98	THR	2.3
14	5A	24	CYS	2.3
20	BA	87	LYS	2.3
28	19	35	LYS	2.3
31	49	11	TYR	2.3
33	69	9	LEU	2.3
31	41	160	VAL	2.3
1	1G	1370	G	2.3
9	8E	71	SER	2.3
26	14	2000	G	2.3
4	32	14	ARG	2.3
32	59	124	GLU	2.3
20	BI	11	SER	2.3
36	35	25	SER	2.3
16	7A	27	LYS	2.3
33	69	45	LYS	2.3
1	1G	1348	U	2.3
2	12	33	TYR	2.3
2	12	70	PHE	2.3
4	32	133	VAL	2.3
16	7I	24	ALA	2.3
20	BI	75	ASN	2.3
29	29	52	LEU	2.3
29	21	160	TYR	2.3
40	B8	50	ILE	2.3
19	AA	51	VAL	2.3
40	75	1	MET	2.3
55	Q8	4	MET	2.3
38	55	24	GLN	2.3
4	3E	68	TYR	2.3
26	1H	2113	U	2.3
42	D8	77	ALA	2.3
1	1G	781	A	2.3
31	49	142	PRO	2.3
4	3E	158	ILE	2.3
9	82	83	ARG	2.3
13	4A	91	ARG	2.3
29	21	141	ILE	2.3
31	49	91	ARG	2.3
37	45	49	ALA	2.3
1	1G	1362(A)	C	2.3
12	3A	17	LYS	2.3

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Mol	Chain	Res	Type	RSRZ
28	19	223	GLY	2.3
20	BA	20	LEU	2.3
28	11	206	LEU	2.3
28	19	16	MET	2.3
28	19	215	LEU	2.3
29	29	195	LEU	2.3
4	3E	204	ILE	2.3
45	C5	75	ILE	2.3
46	D5	153	SER	2.3
3	22	207	VAL	2.3
10	1I	72	VAL	2.3
29	29	135	HIS	2.3
33	69	112	LYS	2.3
50	H5	21	ALA	2.3
26	14	574	C	2.3
28	11	254	THR	2.3
15	6I	56	LEU	2.3
17	8A	43	LEU	2.3
49	G5	61	LEU	2.3
48	F5	33	LYS	2.3
19	AA	14	HIS	2.3
26	1H	2062	A	2.3
42	95	73	SER	2.3
16	7I	71	ARG	2.2
20	BI	64	ASP	2.2
37	45	3	MET	2.2
37	45	75	THR	2.2
41	C8	18	LEU	2.2
49	G5	44	LEU	2.2
11	2I	123	LYS	2.2
12	3I	46	LYS	2.2
28	19	184	LYS	2.2
42	D8	36	PRO	2.2
42	D8	84	LYS	2.2
55	Q8	23	VAL	2.2
35	25	74	GLY	2.2
5	42	18	ARG	2.2
26	14	2001	A	2.2
38	98	68	ARG	2.2
7	62	104	LEU	2.2
51	I5	41	PRO	2.2
1	13	136	C	2.2

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Mol	Chain	Res	Type	RSRZ
5	42	122	GLU	2.2
26	1H	34	C	2.2
26	14	2690	C	2.2
46	D5	99	TYR	2.2
53	O8	24	GLU	2.2
53	O8	51	GLU	2.2
34	15	104	LYS	2.2
55	Q8	12	LYS	2.2
26	1H	1762	A	2.2
26	14	983	A	2.2
41	C8	90	VAL	2.2
41	85	44	ASN	2.2
21	1F	19	GLY	2.2
26	14	2002	G	2.2
26	14	2506	U	2.2
4	3E	61	LYS	2.2
28	19	258	LYS	2.2
38	98	40	LYS	2.2
37	88	17	LEU	2.2
5	4E	19	MET	2.2
31	49	70	VAL	2.2
35	25	2	ILE	2.2
47	E5	60	PHE	2.2
2	1E	15	VAL	2.2
8	72	54	ASP	2.2
10	1A	53	PRO	2.2
20	BI	25	ARG	2.2
47	I8	26	TYR	2.2
1	1G	1150	U	2.2
32	59	158	HIS	2.2
1	13	324	G	2.2
26	14	2110	G	2.2
42	D8	73	SER	2.2
1	1G	1149	C	2.2
47	E5	25	ARG	2.2
19	AA	52	TYR	2.2
32	51	85	LYS	2.2
32	59	91	GLY	2.2
29	21	48	GLN	2.2
47	E5	43	THR	2.2
26	14	2900	A	2.2
31	49	82	LEU	2.2

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Mol	Chain	Res	Type	RSRZ
44	B5	92	LEU	2.2
1	1G	940	C	2.2
11	2A	50	TYR	2.2
36	35	44	GLY	2.2
20	BA	59	ALA	2.2
36	35	106	LEU	2.2
2	12	41	ILE	2.2
26	14	2712	U	2.2
54	L5	2	LYS	2.2
13	4A	8	GLU	2.2
30	39	89	VAL	2.2
29	21	5	LEU	2.2
36	35	81	GLN	2.2
1	13	731	G	2.2
16	7A	28	ARG	2.2
30	39	51	THR	2.2
31	41	33	ARG	2.2
20	BI	55	ILE	2.2
29	29	25	VAL	2.2
1	13	1287	A	2.2
41	85	48	ALA	2.2
9	8E	11	LYS	2.2
11	2A	120	ARG	2.2
16	7A	35	LYS	2.2
43	A5	85	VAL	2.2
1	13	1356	G	2.2
1	1G	769	G	2.2
10	1A	7	LYS	2.2
12	3I	89	ARG	2.2
13	4A	27	LYS	2.2
17	8A	100	LYS	2.2
26	14	2689	U	2.2
36	35	17	LYS	2.2
37	45	18	LYS	2.2
41	85	46	ALA	2.2
32	59	123	PHE	2.2
5	4E	82	VAL	2.2
8	72	132	GLU	2.2
37	45	47	ILE	2.2
9	82	93	ARG	2.2
14	5I	34	TYR	2.2
20	BI	30	LYS	2.2

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Mol	Chain	Res	Type	RSRZ
45	C5	2	ARG	2.2
20	BA	36	LEU	2.2
45	C5	26	LYS	2.2
48	J8	25	LYS	2.2
47	I8	85	ALA	2.2
1	1G	1511	G	2.2
36	78	68	GLN	2.2
39	65	29	PHE	2.2
37	88	38	GLU	2.2
41	85	8	VAL	2.2
1	1G	1288	A	2.2
12	3A	89	ARG	2.2
26	14	575	A	2.2
26	14	2126	A	2.2
4	32	54	TYR	2.1
5	42	133	TYR	2.1
45	C5	31	LEU	2.1
48	F5	71	TYR	2.1
26	14	2477	C	2.1
1	1G	723	U	2.1
37	45	113	GLN	2.1
29	21	171	GLU	2.1
46	D5	167	PRO	2.1
1	1G	1520	G	2.1
1	1G	1529	G	2.1
15	6I	57	LEU	2.1
15	6A	32	LEU	2.1
28	19	61	LEU	2.1
35	25	34	THR	2.1
42	95	71	LEU	2.1
15	6A	15	PHE	2.1
41	C8	117	GLN	2.1
4	32	73	ARG	2.1
10	1A	72	VAL	2.1
36	35	23	PRO	2.1
37	45	127	ILE	2.1
39	65	15	ARG	2.1
45	C5	50	ARG	2.1
55	Q8	56	GLU	2.1
7	62	31	MET	2.1
38	98	1	MET	2.1
55	M5	20	GLY	2.1

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Mol	Chain	Res	Type	RSRZ
5	4E	119	LEU	2.1
33	61	116	LEU	2.1
47	E5	37	LEU	2.1
36	78	57	THR	2.1
48	F5	60	PHE	2.1
11	2I	120	ARG	2.1
12	3A	9	GLN	2.1
24	1L	31	A	2.1
36	35	55	ARG	2.1
41	C8	15	LYS	2.1
29	29	126	PRO	2.1
32	59	55	PRO	2.1
38	55	6	SER	2.1
1	1G	980	C	2.1
4	3E	97	LEU	2.1
23	2L	1	C	2.1
4	32	4	TYR	2.1
30	31	97	TYR	2.1
47	I8	77	ARG	2.1
55	M5	17	THR	2.1
19	AA	11	VAL	2.1
20	BA	60	GLU	2.1
28	11	18	VAL	2.1
28	19	49	ILE	2.1
28	19	64	ILE	2.1
30	39	88	VAL	2.1
36	78	71	VAL	2.1
48	J8	13	ILE	2.1
48	J8	49	VAL	2.1
8	7E	59	LEU	2.1
26	14	859	G	2.1
26	14	2052	G	2.1
1	1G	1020	U	2.1
7	62	29	LYS	2.1
11	2I	122	LYS	2.1
11	2A	75	TYR	2.1
26	14	2723	C	2.1
30	39	44	ARG	2.1
50	H5	18	ASP	2.1
28	19	214	TRP	2.1
55	M5	31	HIS	2.1
20	BA	9	ASN	2.1

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Mol	Chain	Res	Type	RSRZ
51	I5	34	GLU	2.1
12	3A	70	ILE	2.1
32	59	71	LEU	2.1
8	72	21	LYS	2.1
32	51	170	ARG	2.1
47	E5	20	ARG	2.1
55	M5	36	LYS	2.1
1	13	1531	A	2.1
1	13	107	G	2.1
24	3L	6	G	2.1
26	14	2120	G	2.1
8	7E	24	THR	2.1
26	14	280	C	2.1
4	3E	76	ARG	2.1
4	32	49	ARG	2.1
21	1F	12	LYS	2.1
21	1B	7	ARG	2.1
32	59	30	LYS	2.1
35	25	70	LYS	2.1
28	11	53	PHE	2.1
7	62	7	ALA	2.1
7	62	40	ALA	2.1
38	98	21	TYR	2.1
8	7E	134	ILE	2.1
32	59	148	ILE	2.1
4	3E	122	ARG	2.1
5	4E	31	LEU	2.1
12	3A	72	GLY	2.1
26	14	530	G	2.1
28	19	161	THR	2.1
29	21	49	LEU	2.1
44	F8	2	LYS	2.1
52	J5	3	LYS	2.1
8	72	27	PRO	2.1
47	E5	26	TYR	2.1
28	19	181	GLU	2.1
28	11	35	LYS	2.1
28	19	221	VAL	2.1
44	B5	52	VAL	2.1
5	4E	24	ARG	2.1
29	29	124	GLY	2.1
31	49	177	GLY	2.1

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Mol	Chain	Res	Type	RSRZ
36	78	106	LEU	2.1
37	45	56	ARG	2.1
48	F5	23	LYS	2.1
15	6A	66	LEU	2.1
26	14	1301	A	2.1
1	13	625	G	2.1
13	4A	67	GLU	2.1
32	59	81	GLU	2.1
39	65	43	GLU	2.1
45	C5	33	LYS	2.1
47	E5	46	LYS	2.1
50	H5	30	ARG	2.1
20	BA	24	LEU	2.1
47	I8	54	GLY	2.1
55	M5	45	GLY	2.1
34	15	43	THR	2.1
1	13	1350	A	2.1
4	3E	80	GLU	2.1
20	BI	34	LYS	2.1
30	39	49	ALA	2.1
42	D8	78	LYS	2.1
43	A5	16	LYS	2.1
9	8E	17	VAL	2.1
55	Q8	42	ARG	2.1
5	4E	118	ILE	2.1
24	1L	56	C	2.1
4	3E	120	LEU	2.1
32	59	33	LEU	2.1
34	15	113	GLY	2.1
34	15	120	LEU	2.1
1	1G	1235	U	2.1
14	5A	46	GLU	2.1
14	5I	18	VAL	2.1
1	13	623	C	2.1
8	7E	25	ASP	2.1
18	9A	43	PHE	2.0
39	A8	112	PHE	2.0
42	D8	76	LYS	2.0
52	N8	15	ARG	2.0
26	14	352	G	2.0
5	42	17	ALA	2.0
8	72	16	ALA	2.0

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Mol	Chain	Res	Type	RSRZ
37	45	132	VAL	2.0
48	F5	4	VAL	2.0
13	4I	6	GLY	2.0
4	3E	84	LYS	2.0
30	39	75	HIS	2.0
36	78	79	ARG	2.0
42	D8	80	GLN	2.0
51	I5	49	PHE	2.0
7	62	33	ASP	2.0
47	I8	56	ASP	2.0
1	13	110	C	2.0
17	8I	24	GLU	2.0
12	3A	27	LEU	2.0
29	29	138	PRO	2.0
32	59	108	GLY	2.0
34	58	107	LEU	2.0
26	14	573	G	2.0
37	88	33	GLY	2.0
37	88	130	LYS	2.0
16	7A	8	ARG	2.0
40	75	112	ARG	2.0
44	B5	28	PHE	2.0
47	I8	44	ARG	2.0
9	8E	12	GLU	2.0
1	13	608	A	2.0
4	3E	140	VAL	2.0
14	5I	21	TYR	2.0
28	19	51	VAL	2.0
48	F5	74	VAL	2.0
1	13	1344	C	2.0
11	2A	71	LYS	2.0
16	7I	73	LEU	2.0
16	7I	63	GLY	2.0
28	19	17	THR	2.0
29	29	148	GLY	2.0
35	25	67	LYS	2.0
28	19	253	GLN	2.0
2	12	71	VAL	2.0
10	1I	73	ASP	2.0
39	65	27	SER	2.0
12	3A	97	ARG	2.0
19	AA	37	ARG	2.0

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Mol	Chain	Res	Type	RSRZ
19	AA	62	ILE	2.0
28	19	204	ILE	2.0
29	29	118	LYS	2.0
38	55	44	LEU	2.0
39	65	26	LEU	2.0
48	J8	23	LYS	2.0
44	B5	60	ARG	2.0
50	H5	26	LEU	2.0
52	N8	25	LEU	2.0
29	29	193	GLY	2.0
1	1G	1066	C	2.0
26	14	1625	C	2.0
29	29	147	PRO	2.0
38	98	39	PRO	2.0
37	45	45	GLN	2.0
5	42	12	LEU	2.0
12	3I	7	ILE	2.0
15	6I	32	LEU	2.0
36	35	148	LEU	2.0
37	88	32	TYR	2.0
31	49	127	GLY	2.0
51	M8	31	ILE	2.0
1	13	1361	G	2.0
1	1G	60	A	2.0
1	1G	815	A	2.0
9	8E	103	THR	2.0
22	1K	3	C	2.0
24	1L	2	C	2.0
38	55	40	LYS	2.0
4	3E	108	LEU	2.0
7	6E	154	TYR	2.0
37	88	132	VAL	2.0
55	M5	60	LEU	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
22	PSU	1K	39	20/21	0.94	0.17	-	96,108,116,119	0
23	4SU	2K	8	20/21	0.94	0.21	-	100,108,123,123	0
22	PSU	1K	32	20/21	0.89	0.32	-	113,119,123,123	0
23	OMC	2L	33	21/22	0.96	0.20	-	124,127,132,137	0
22	PSU	1K	55	20/21	0.78	0.11	-	153,172,191,191	0
23	OMC	2K	33	21/22	0.98	0.32	-	88,94,101,106	0
23	4SU	2L	8	20/21	0.88	0.14	-	128,140,143,147	0
23	PSU	2K	56	20/21	0.96	0.14	-	109,113,120,127	0
22	4SU	1K	8	20/21	0.80	0.14	-	184,192,200,205	0
23	5MU	2L	55	21/22	0.95	0.11	-	138,142,147,151	0
22	7MG	1K	46	24/25	0.87	0.12	-	175,197,203,209	0
23	7MG	2K	47	24/25	0.95	0.17	-	110,120,128,137	0
23	PSU	2L	56	20/21	0.90	0.09	-	128,135,142,148	0
23	5MU	2K	55	21/22	0.97	0.15	-	99,116,125,136	0
22	5MU	1K	54	21/22	0.93	0.10	-	142,157,168,172	0
23	7MG	2L	47	24/25	0.95	0.12	-	140,151,156,161	0
22	MIA	1K	37	29/30	0.94	0.30	-	89,100,118,123	0

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
56	MG	13	1641	1/1	0.76	0.35	34.96	88,88,88,88	0
56	MG	1H	3173	1/1	0.95	0.36	28.63	98,98,98,98	0
56	MG	1H	3277	1/1	0.90	0.47	18.53	109,109,109,109	0
56	MG	13	1621	1/1	0.87	0.39	16.07	115,115,115,115	0
56	MG	1H	3025	1/1	0.95	0.48	15.77	78,78,78,78	0
56	MG	14	3217	1/1	0.90	0.80	15.55	82,82,82,82	0
56	MG	2K	105	1/1	0.80	0.38	14.78	107,107,107,107	0
56	MG	1H	3084	1/1	0.98	0.31	14.73	61,61,61,61	0
56	MG	1H	3037	1/1	0.97	0.37	14.32	76,76,76,76	0
56	MG	1H	3245	1/1	0.89	0.38	13.50	96,96,96,96	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
56	MG	1H	3069	1/1	0.96	0.37	11.34	89,89,89,89	0
56	MG	1H	3298	1/1	0.85	0.33	11.24	92,92,92,92	0
56	MG	1H	3304	1/1	0.90	0.39	11.21	114,114,114,114	0
56	MG	1H	3011	1/1	0.95	0.51	10.37	91,91,91,91	0
56	MG	16	206	1/1	0.86	0.25	9.60	102,102,102,102	0
56	MG	14	3274	1/1	0.72	0.25	9.53	122,122,122,122	0
56	MG	13	1686	1/1	0.92	0.29	8.88	134,134,134,134	0
56	MG	14	3169	1/1	0.82	0.29	8.48	93,93,93,93	0
56	MG	1H	3255	1/1	0.63	0.45	8.39	117,117,117,117	0
56	MG	1H	3056	1/1	0.92	0.30	8.23	72,72,72,72	0
56	MG	1H	3234	1/1	0.98	0.29	8.18	91,91,91,91	0
56	MG	13	1689	1/1	0.74	0.25	8.16	98,98,98,98	0
56	MG	14	3170	1/1	0.75	0.60	8.02	77,77,77,77	0
56	MG	1H	3013	1/1	0.80	0.41	8.01	104,104,104,104	0
56	MG	1H	3222	1/1	0.87	0.30	7.88	79,79,79,79	0
56	MG	14	3199	1/1	0.96	0.35	7.78	126,126,126,126	0
56	MG	1H	3170	1/1	0.95	0.33	7.71	100,100,100,100	0
56	MG	16	207	1/1	0.90	0.37	7.65	84,84,84,84	0
56	MG	1G	1660	1/1	0.93	0.43	7.50	97,97,97,97	0
56	MG	14	3055	1/1	0.91	0.42	7.25	85,85,85,85	0
56	MG	14	3233	1/1	0.88	0.30	7.09	118,118,118,118	0
56	MG	1H	3160	1/1	0.93	0.31	7.06	84,84,84,84	0
56	MG	1G	1614	1/1	0.93	0.25	7.03	140,140,140,140	0
56	MG	1H	3258	1/1	0.85	0.31	6.54	93,93,93,93	0
56	MG	1H	3110	1/1	0.97	0.31	6.50	67,67,67,67	0
56	MG	1H	3186	1/1	0.70	0.35	6.20	93,93,93,93	0
56	MG	14	3216	1/1	0.95	0.43	5.98	111,111,111,111	0
56	MG	L8	101	1/1	0.79	0.43	5.82	85,85,85,85	0
56	MG	14	3306	1/1	0.90	0.30	5.76	93,93,93,93	0
56	MG	14	3091	1/1	0.93	0.29	5.74	87,87,87,87	0
56	MG	14	3272	1/1	0.71	0.32	5.72	95,95,95,95	0
56	MG	1H	3063	1/1	0.98	0.28	5.57	83,83,83,83	0
56	MG	14	3195	1/1	0.86	0.38	5.47	100,100,100,100	0
56	MG	1H	3116	1/1	0.80	0.29	5.41	70,70,70,70	0
56	MG	14	3253	1/1	0.83	0.29	5.36	82,82,82,82	0
56	MG	14	3227	1/1	0.70	0.33	5.22	97,97,97,97	0
56	MG	14	3211	1/1	0.82	0.40	5.14	93,93,93,93	0
56	MG	1H	3235	1/1	0.95	0.25	4.93	75,75,75,75	0
56	MG	14	3202	1/1	0.84	0.26	4.91	100,100,100,100	0
56	MG	1H	3254	1/1	0.82	0.37	4.86	80,80,80,80	0
56	MG	39	303	1/1	0.66	0.67	4.80	87,87,87,87	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
56	MG	1H	3178	1/1	0.86	0.27	4.68	76,76,76,76	0
56	MG	14	3194	1/1	0.97	0.23	4.27	97,97,97,97	0
56	MG	1H	3219	1/1	0.83	0.39	4.23	65,65,65,65	0
56	MG	1H	3089	1/1	0.97	0.27	3.98	58,58,58,58	0
56	MG	1H	3004	1/1	0.99	0.30	3.86	68,68,68,68	0
56	MG	14	3205	1/1	0.97	0.45	3.80	72,72,72,72	0
56	MG	13	1664	1/1	0.65	0.28	3.68	103,103,103,103	0
56	MG	1H	3081	1/1	0.84	0.25	3.53	77,77,77,77	0
56	MG	1H	3332	1/1	0.63	0.42	3.50	90,90,90,90	0
56	MG	13	1706	1/1	0.88	0.91	3.50	103,103,103,103	0
56	MG	29	301	1/1	0.74	0.64	3.50	101,101,101,101	0
56	MG	14	3209	1/1	0.85	0.33	3.49	71,71,71,71	0
56	MG	13	1602	1/1	0.94	0.31	3.43	87,87,87,87	0
56	MG	1G	1659	1/1	0.81	0.30	3.35	102,102,102,102	0
56	MG	14	3382	1/1	0.97	0.29	3.28	86,86,86,86	0
56	MG	13	1652	1/1	0.88	0.23	3.20	98,98,98,98	0
56	MG	1G	1645	1/1	0.68	0.31	3.18	113,113,113,113	0
56	MG	1H	3052	1/1	0.95	0.30	3.15	80,80,80,80	0
56	MG	14	3308	1/1	0.85	0.21	3.09	106,106,106,106	0
56	MG	14	3264	1/1	0.76	0.30	2.72	89,89,89,89	0
56	MG	1J	207	1/1	0.87	0.27	2.71	147,147,147,147	0
56	MG	1H	3260	1/1	0.86	0.29	2.57	76,76,76,76	0
56	MG	14	3266	1/1	0.86	0.38	2.55	80,80,80,80	0
56	MG	1H	3139	1/1	0.96	0.40	2.49	79,79,79,79	0
56	MG	14	3007	1/1	0.95	0.23	2.47	94,94,94,94	0
56	MG	1H	3240	1/1	0.96	0.25	2.43	94,94,94,94	0
56	MG	13	1680	1/1	0.88	0.19	2.33	124,124,124,124	0
56	MG	14	3248	1/1	0.80	0.26	2.33	85,85,85,85	0
56	MG	1G	1664	1/1	0.67	0.27	2.30	116,116,116,116	0
56	MG	13	1666	1/1	0.71	0.32	2.29	105,105,105,105	0
56	MG	16	209	1/1	0.96	0.21	2.25	88,88,88,88	0
56	MG	14	3060	1/1	0.95	0.27	2.23	83,83,83,83	0
56	MG	14	3146	1/1	0.93	0.19	2.22	128,128,128,128	0
56	MG	1G	1627	1/1	0.97	0.23	2.20	116,116,116,116	0
57	ZN	32	302	1/1	0.99	0.39	2.09	131,131,131,131	0
56	MG	14	3254	1/1	0.81	0.17	2.09	87,87,87,87	0
56	MG	14	3165	1/1	0.96	0.33	1.96	70,70,70,70	0
56	MG	13	1712	1/1	0.97	0.20	1.94	99,99,99,99	0
56	MG	14	3073	1/1	0.96	0.26	1.93	74,74,74,74	0
56	MG	1H	3230	1/1	0.79	0.24	1.89	91,91,91,91	0
56	MG	13	1603	1/1	0.94	0.21	1.86	97,97,97,97	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
56	MG	1H	3205	1/1	0.99	0.23	1.64	64,64,64,64	0
56	MG	14	3049	1/1	0.95	0.25	1.59	87,87,87,87	0
56	MG	16	214	1/1	0.78	0.22	1.55	109,109,109,109	0
56	MG	13	1721	1/1	0.77	0.63	1.53	101,101,101,101	0
56	MG	14	3150	1/1	0.66	0.33	1.50	80,80,80,80	0
56	MG	13	1604	1/1	0.98	0.27	1.50	108,108,108,108	0
56	MG	1H	3064	1/1	0.96	0.27	1.45	76,76,76,76	0
56	MG	41	202	1/1	0.79	0.29	1.40	96,96,96,96	0
56	MG	1H	3088	1/1	0.97	0.33	1.24	52,52,52,52	0
56	MG	I8	101	1/1	0.81	0.27	1.24	84,84,84,84	0
56	MG	14	3220	1/1	0.96	0.32	1.19	94,94,94,94	0
56	MG	1H	3216	1/1	0.92	0.22	1.05	77,77,77,77	0
56	MG	14	3262	1/1	0.54	0.29	1.02	81,81,81,81	0
56	MG	1H	3288	1/1	0.84	0.28	0.99	87,87,87,87	0
56	MG	14	3267	1/1	0.83	0.16	0.93	110,110,110,110	0
56	MG	13	1648	1/1	0.96	0.26	0.92	76,76,76,76	0
56	MG	14	3119	1/1	0.98	0.32	0.89	66,66,66,66	0
56	MG	1H	3001	1/1	0.98	0.29	0.89	64,64,64,64	0
56	MG	14	3268	1/1	0.89	0.19	0.84	100,100,100,100	0
56	MG	14	3058	1/1	0.85	0.15	0.84	128,128,128,128	0
56	MG	1H	3172	1/1	0.97	0.26	0.83	98,98,98,98	0
57	ZN	G8	202	1/1	0.87	0.31	0.82	192,192,192,192	0
56	MG	16	205	1/1	0.53	0.20	0.80	91,91,91,91	0
56	MG	13	1601	1/1	0.98	0.27	0.79	74,74,74,74	0
57	ZN	3E	301	1/1	0.98	0.41	0.79	115,115,115,115	0
56	MG	1H	3273	1/1	0.68	0.25	0.76	84,84,84,84	0
56	MG	14	3260	1/1	0.63	0.17	0.75	119,119,119,119	0
56	MG	14	3110	1/1	0.98	0.23	0.71	61,61,61,61	0
56	MG	13	1612	1/1	0.96	0.27	0.66	89,89,89,89	0
56	MG	13	1618	1/1	0.90	0.26	0.66	86,86,86,86	0
56	MG	14	3411	1/1	0.84	0.30	0.65	104,104,104,104	0
56	MG	14	3053	1/1	0.92	0.17	0.62	98,98,98,98	0
56	MG	1H	3241	1/1	0.91	0.21	0.59	80,80,80,80	0
56	MG	14	3128	1/1	0.97	0.21	0.57	97,97,97,97	0
56	MG	1H	3252	1/1	0.93	0.21	0.55	69,69,69,69	0
56	MG	1H	3077	1/1	0.90	0.25	0.53	67,67,67,67	0
56	MG	13	1625	1/1	0.95	0.30	0.52	77,77,77,77	0
56	MG	1H	3076	1/1	0.99	0.26	0.43	73,73,73,73	0
56	MG	14	3014	1/1	0.99	0.27	0.43	73,73,73,73	0
56	MG	14	3256	1/1	0.83	0.19	0.42	97,97,97,97	0
56	MG	14	3118	1/1	0.98	0.24	0.42	71,71,71,71	0
56	MG	14	3098	1/1	0.92	0.28	0.42	74,74,74,74	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
56	MG	14	3121	1/1	0.94	0.23	0.38	84,84,84,84	0
56	MG	13	1752	1/1	0.89	0.51	0.38	92,92,92,92	0
56	MG	14	3106	1/1	0.94	0.20	0.37	89,89,89,89	0
56	MG	1G	1623	1/1	0.83	0.24	0.35	122,122,122,122	0
56	MG	13	1607	1/1	0.89	0.19	0.34	78,78,78,78	0
56	MG	13	1682	1/1	0.83	0.28	0.31	116,116,116,116	0
56	MG	16	203	1/1	0.97	0.18	0.26	107,107,107,107	0
56	MG	1G	1655	1/1	0.91	0.21	0.25	111,111,111,111	0
56	MG	1H	3244	1/1	0.96	0.20	0.23	90,90,90,90	0
56	MG	14	3278	1/1	0.86	0.26	0.21	85,85,85,85	0
56	MG	14	3154	1/1	0.91	0.23	0.19	86,86,86,86	0
56	MG	1H	3203	1/1	0.94	0.35	-0.07	73,73,73,73	0
56	MG	1H	3184	1/1	0.90	0.19	-0.09	107,107,107,107	0
56	MG	13	1667	1/1	0.92	0.30	-0.12	147,147,147,147	0
56	MG	14	3040	1/1	0.96	0.20	-0.12	79,79,79,79	0
56	MG	J8	101	1/1	0.94	0.26	-0.13	78,78,78,78	0
56	MG	1G	1607	1/1	0.93	0.28	-0.20	98,98,98,98	0
56	MG	14	3252	1/1	0.74	0.15	-0.20	105,105,105,105	0
56	MG	13	1620	1/1	0.94	0.28	-0.21	119,119,119,119	0
56	MG	85	201	1/1	0.95	0.28	-0.22	93,93,93,93	0
56	MG	14	3062	1/1	0.97	0.29	-0.23	85,85,85,85	0
56	MG	14	3052	1/1	0.94	0.21	-0.25	110,110,110,110	0
56	MG	14	3138	1/1	0.95	0.20	-0.28	73,73,73,73	0
56	MG	1H	3155	1/1	0.87	0.17	-0.29	78,78,78,78	0
56	MG	1H	3315	1/1	0.94	0.39	-0.32	105,105,105,105	0
56	MG	1H	3130	1/1	0.98	0.24	-0.33	78,78,78,78	0
56	MG	1G	1628	1/1	0.83	0.21	-0.34	118,118,118,118	0
56	MG	14	3219	1/1	0.95	0.20	-0.39	102,102,102,102	0
56	MG	14	3107	1/1	0.92	0.21	-0.42	80,80,80,80	0
56	MG	14	3103	1/1	0.78	0.38	-0.44	90,90,90,90	0
56	MG	14	3238	1/1	0.97	0.16	-0.46	110,110,110,110	0
56	MG	1H	3470	1/1	0.93	0.23	-0.47	99,99,99,99	0
56	MG	1G	1635	1/1	0.87	0.39	-0.48	118,118,118,118	0
56	MG	1H	3085	1/1	0.98	0.22	-0.49	62,62,62,62	0
56	MG	14	3008	1/1	0.98	0.25	-0.50	78,78,78,78	0
56	MG	13	1728	1/1	0.94	0.13	-0.51	131,131,131,131	0
56	MG	14	3244	1/1	0.97	0.23	-0.60	99,99,99,99	0
56	MG	13	1675	1/1	0.71	0.19	-0.65	121,121,121,121	0
56	MG	14	3215	1/1	0.99	0.15	-0.65	104,104,104,104	0
56	MG	1G	1644	1/1	0.59	0.19	-0.66	115,115,115,115	0
56	MG	1H	3324	1/1	0.95	0.15	-0.70	107,107,107,107	0
56	MG	13	1611	1/1	0.99	0.18	-0.71	95,95,95,95	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
56	MG	88	201	1/1	0.95	0.22	-0.72	95,95,95,95	0
56	MG	13	1743	1/1	0.77	0.18	-0.76	108,108,108,108	0
56	MG	13	1751	1/1	0.96	0.16	-0.77	94,94,94,94	0
56	MG	1J	204	1/1	0.82	0.12	-0.78	113,113,113,113	0
57	ZN	5I	101	1/1	0.97	0.20	-0.81	114,114,114,114	0
56	MG	14	3079	1/1	0.99	0.23	-0.82	84,84,84,84	0
56	MG	13	1608	1/1	0.97	0.20	-0.85	91,91,91,91	0
56	MG	14	3051	1/1	0.98	0.21	-0.85	100,100,100,100	0
56	MG	3I	201	1/1	0.95	0.19	-0.86	81,81,81,81	0
56	MG	14	3084	1/1	0.96	0.22	-0.86	78,78,78,78	0
56	MG	1H	3009	1/1	0.90	0.20	-0.87	69,69,69,69	0
56	MG	14	3171	1/1	0.95	0.19	-0.87	73,73,73,73	0
56	MG	11	301	1/1	0.92	0.31	-0.88	62,62,62,62	0
56	MG	14	3137	1/1	0.78	0.18	-0.88	80,80,80,80	0
56	MG	14	3009	1/1	0.96	0.24	-0.93	65,65,65,65	0
56	MG	14	3115	1/1	0.95	0.18	-0.95	100,100,100,100	0
56	MG	1G	1653	1/1	0.86	0.11	-0.96	106,106,106,106	0
56	MG	14	3425	1/1	0.94	0.28	-0.97	92,92,92,92	0
56	MG	2K	101	1/1	0.97	0.22	-0.99	76,76,76,76	0
56	MG	13	1671	1/1	0.94	0.24	-1.00	102,102,102,102	0
56	MG	14	3070	1/1	0.96	0.19	-1.01	66,66,66,66	0
56	MG	1H	3421	1/1	0.94	0.18	-1.06	62,62,62,62	0
56	MG	2L	101	1/1	0.97	0.18	-1.06	98,98,98,98	0
56	MG	1H	3053	1/1	0.94	0.17	-1.12	92,92,92,92	0
56	MG	1H	3101	1/1	0.92	0.21	-1.14	49,49,49,49	0
56	MG	1G	1667	1/1	0.67	0.24	-1.15	119,119,119,119	0
56	MG	39	301	1/1	0.93	0.19	-1.16	111,111,111,111	0
56	MG	14	3208	1/1	0.85	0.10	-1.16	84,84,84,84	0
56	MG	1H	3020	1/1	0.93	0.15	-1.17	90,90,90,90	0
56	MG	1H	3054	1/1	0.96	0.19	-1.17	71,71,71,71	0
56	MG	1G	1680	1/1	0.82	0.11	-1.20	140,140,140,140	0
56	MG	45	201	1/1	0.86	0.18	-1.22	113,113,113,113	0
56	MG	13	1668	1/1	0.80	0.19	-1.25	94,94,94,94	0
56	MG	14	3120	1/1	0.98	0.17	-1.25	80,80,80,80	0
56	MG	1H	3456	1/1	0.94	0.18	-1.31	79,79,79,79	0
56	MG	1H	3166	1/1	0.93	0.17	-1.32	73,73,73,73	0
56	MG	14	3221	1/1	0.88	0.20	-1.34	87,87,87,87	0
57	ZN	5A	101	1/1	0.95	0.13	-1.36	172,172,172,172	0
56	MG	13	1655	1/1	0.96	0.14	-1.36	89,89,89,89	0
56	MG	13	1660	1/1	0.93	0.14	-1.39	94,94,94,94	0
56	MG	1H	3209	1/1	0.90	0.15	-1.41	82,82,82,82	0
56	MG	1H	3006	1/1	0.98	0.17	-1.42	91,91,91,91	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
56	MG	1G	1688	1/1	0.91	0.13	-1.44	117,117,117,117	0
56	MG	41	201	1/1	0.95	0.16	-1.44	89,89,89,89	0
56	MG	49	201	1/1	0.88	0.10	-1.44	143,143,143,143	0
56	MG	14	3041	1/1	0.94	0.16	-1.46	80,80,80,80	0
56	MG	14	3310	1/1	0.97	0.20	-1.47	96,96,96,96	0
56	MG	14	3076	1/1	0.90	0.11	-1.48	93,93,93,93	0
56	MG	1J	201	1/1	0.95	0.15	-1.48	123,123,123,123	0
56	MG	1H	3368	1/1	0.99	0.17	-1.50	79,79,79,79	0
56	MG	1G	1601	1/1	0.98	0.19	-1.57	109,109,109,109	0
56	MG	1H	3003	1/1	0.93	0.21	-1.58	75,75,75,75	0
56	MG	1H	3040	1/1	0.99	0.21	-1.61	78,78,78,78	0
56	MG	14	3246	1/1	0.92	0.21	-1.65	75,75,75,75	0
56	MG	1H	3067	1/1	0.94	0.16	-1.65	84,84,84,84	0
56	MG	14	3379	1/1	0.90	0.09	-1.66	77,77,77,77	0
56	MG	14	3198	1/1	0.97	0.11	-1.67	88,88,88,88	0
56	MG	14	3072	1/1	0.93	0.18	-1.67	90,90,90,90	0
56	MG	21	302	1/1	0.96	0.16	-1.67	82,82,82,82	0
56	MG	14	3016	1/1	0.99	0.24	-1.69	71,71,71,71	0
56	MG	14	3082	1/1	0.94	0.22	-1.70	71,71,71,71	0
56	MG	1H	3148	1/1	0.65	0.18	-1.76	79,79,79,79	0
56	MG	1H	3393	1/1	0.98	0.17	-1.76	60,60,60,60	0
56	MG	1H	3361	1/1	0.98	0.16	-1.80	56,56,56,56	0
56	MG	13	1723	1/1	0.98	0.14	-1.84	105,105,105,105	0
56	MG	14	3332	1/1	0.99	0.12	-1.85	72,72,72,72	0
56	MG	1H	3180	1/1	0.96	0.17	-1.89	84,84,84,84	0
56	MG	14	3095	1/1	0.97	0.20	-1.89	78,78,78,78	0
56	MG	14	3153	1/1	0.90	0.17	-1.89	70,70,70,70	0
56	MG	1H	3351	1/1	0.98	0.18	-1.91	65,65,65,65	0
56	MG	14	3369	1/1	0.88	0.18	-1.92	105,105,105,105	0
56	MG	14	3300	1/1	0.97	0.15	-1.93	98,98,98,98	0
56	MG	1G	1674	1/1	0.92	0.15	-1.94	111,111,111,111	0
56	MG	1H	3132	1/1	0.80	0.19	-1.94	74,74,74,74	0
56	MG	14	3393	1/1	0.95	0.14	-1.95	80,80,80,80	0
56	MG	14	3390	1/1	0.95	0.18	-1.96	94,94,94,94	0
56	MG	14	3094	1/1	0.92	0.20	-1.96	83,83,83,83	0
56	MG	14	3033	1/1	0.98	0.18	-1.99	66,66,66,66	0
56	MG	14	3189	1/1	0.98	0.12	-2.00	83,83,83,83	0
56	MG	14	3010	1/1	0.98	0.19	-2.04	65,65,65,65	0
56	MG	14	3257	1/1	0.99	0.20	-2.04	84,84,84,84	0
56	MG	1H	3024	1/1	0.94	0.23	-2.06	102,102,102,102	0
56	MG	13	1647	1/1	0.86	0.17	-2.07	93,93,93,93	0
56	MG	13	1679	1/1	0.88	0.19	-2.11	98,98,98,98	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
56	MG	1H	3403	1/1	0.97	0.16	-2.13	73,73,73,73	0
56	MG	1H	3055	1/1	0.83	0.16	-2.16	82,82,82,82	0
56	MG	14	3235	1/1	0.69	0.14	-2.20	83,83,83,83	0
56	MG	1H	3366	1/1	0.90	0.12	-2.24	60,60,60,60	0
56	MG	14	3337	1/1	0.87	0.12	-2.27	76,76,76,76	0
56	MG	14	3365	1/1	0.92	0.10	-2.28	80,80,80,80	0
56	MG	1G	1625	1/1	0.76	0.12	-2.28	131,131,131,131	0
56	MG	1G	1679	1/1	0.88	0.09	-2.30	108,108,108,108	0
56	MG	1H	3353	1/1	0.98	0.11	-2.35	77,77,77,77	0
56	MG	1H	3044	1/1	0.99	0.20	-2.36	68,68,68,68	0
56	MG	1H	3233	1/1	0.98	0.19	-2.38	65,65,65,65	0
56	MG	14	3206	1/1	0.96	0.08	-2.41	83,83,83,83	0
57	ZN	C5	202	1/1	0.89	0.11	-2.42	189,189,189,189	0
56	MG	1H	3145	1/1	0.98	0.20	-2.44	57,57,57,57	0
56	MG	1H	3163	1/1	0.92	0.20	-2.44	88,88,88,88	0
56	MG	1H	3489	1/1	0.85	0.12	-2.46	69,69,69,69	0
56	MG	1H	3096	1/1	0.97	0.17	-2.48	57,57,57,57	0
56	MG	1H	3237	1/1	0.89	0.17	-2.48	70,70,70,70	0
56	MG	1H	3215	1/1	0.95	0.14	-2.49	81,81,81,81	0
56	MG	14	3353	1/1	0.97	0.15	-2.49	91,91,91,91	0
56	MG	1H	3390	1/1	0.94	0.17	-2.52	83,83,83,83	0
56	MG	13	1713	1/1	0.91	0.12	-2.53	119,119,119,119	0
56	MG	1H	3005	1/1	0.92	0.20	-2.55	70,70,70,70	0
56	MG	1H	3087	1/1	0.69	0.18	-2.55	69,69,69,69	0
56	MG	14	3069	1/1	0.95	0.08	-2.62	63,63,63,63	0
56	MG	14	3357	1/1	0.99	0.13	-2.63	80,80,80,80	0
56	MG	14	3387	1/1	0.90	0.10	-2.69	76,76,76,76	0
56	MG	1H	3082	1/1	0.93	0.17	-2.71	77,77,77,77	0
56	MG	13	1650	1/1	0.86	0.15	-2.72	87,87,87,87	0
56	MG	1H	3431	1/1	0.65	0.11	-2.73	163,163,163,163	0
56	MG	13	1729	1/1	0.97	0.12	-2.76	109,109,109,109	0
56	MG	1H	3360	1/1	0.96	0.14	-2.77	58,58,58,58	0
56	MG	1G	1662	1/1	0.95	0.16	-2.83	155,155,155,155	0
56	MG	1H	3124	1/1	0.71	0.19	-2.88	72,72,72,72	0
56	MG	14	3204	1/1	0.83	0.13	-2.89	80,80,80,80	0
56	MG	1G	1682	1/1	0.88	0.09	-2.90	141,141,141,141	0
56	MG	13	1642	1/1	0.95	0.11	-2.91	98,98,98,98	0
56	MG	1H	3364	1/1	0.96	0.12	-2.92	80,80,80,80	0
56	MG	1H	3042	1/1	0.96	0.18	-2.93	53,53,53,53	0
56	MG	14	3044	1/1	0.95	0.14	-2.96	78,78,78,78	0
56	MG	1G	1678	1/1	0.96	0.13	-2.97	108,108,108,108	0
56	MG	13	1732	1/1	0.98	0.09	-3.02	81,81,81,81	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
56	MG	14	3005	1/1	0.95	0.18	-3.06	66,66,66,66	0
56	MG	13	1694	1/1	0.80	0.10	-3.07	121,121,121,121	0
56	MG	14	3192	1/1	0.88	0.14	-3.10	84,84,84,84	0
56	MG	14	3322	1/1	0.97	0.14	-3.17	90,90,90,90	0
56	MG	1H	3388	1/1	0.97	0.14	-3.22	65,65,65,65	0
56	MG	1H	3356	1/1	0.97	0.10	-3.30	58,58,58,58	0
56	MG	1H	3159	1/1	0.94	0.17	-3.33	77,77,77,77	0
56	MG	1H	3211	1/1	0.91	0.14	-3.34	76,76,76,76	0
56	MG	1H	3120	1/1	0.97	0.13	-3.35	80,80,80,80	0
56	MG	1H	3125	1/1	0.98	0.16	-3.35	68,68,68,68	0
56	MG	1H	3113	1/1	0.94	0.14	-3.36	64,64,64,64	0
56	MG	13	1643	1/1	0.94	0.09	-3.41	78,78,78,78	0
56	MG	1G	1609	1/1	0.91	0.12	-3.43	97,97,97,97	0
56	MG	13	1638	1/1	0.97	0.09	-3.44	115,115,115,115	0
56	MG	1G	1608	1/1	0.98	0.14	-3.47	98,98,98,98	0
56	MG	1H	3379	1/1	0.98	0.12	-3.47	74,74,74,74	0
56	MG	14	3358	1/1	0.98	0.09	-3.50	85,85,85,85	0
56	MG	14	3111	1/1	0.97	0.17	-3.51	72,72,72,72	0
56	MG	14	3336	1/1	0.95	0.14	-3.61	88,88,88,88	0
56	MG	14	3344	1/1	0.92	0.11	-3.66	82,82,82,82	0
56	MG	13	1662	1/1	0.91	0.11	-3.67	94,94,94,94	0
56	MG	14	3061	1/1	0.98	0.15	-3.68	79,79,79,79	0
56	MG	1G	1618	1/1	0.94	0.12	-3.68	104,104,104,104	0
56	MG	13	1738	1/1	0.98	0.13	-3.75	77,77,77,77	0
56	MG	13	1695	1/1	0.82	0.15	-3.79	112,112,112,112	0
56	MG	1G	1622	1/1	0.93	0.12	-3.79	89,89,89,89	0
56	MG	14	3047	1/1	0.91	0.14	-3.84	66,66,66,66	0
56	MG	1H	3369	1/1	0.97	0.15	-3.84	92,92,92,92	0
56	MG	1H	3471	1/1	0.88	0.11	-3.86	128,128,128,128	0
56	MG	1H	3385	1/1	0.95	0.14	-3.89	75,75,75,75	0
56	MG	14	3236	1/1	0.71	0.10	-3.95	99,99,99,99	0
56	MG	13	1617	1/1	0.96	0.08	-3.96	106,106,106,106	0
56	MG	1H	3350	1/1	0.96	0.14	-3.98	64,64,64,64	0
56	MG	1H	3362	1/1	0.99	0.14	-4.02	61,61,61,61	0
56	MG	1H	3251	1/1	0.72	0.15	-4.03	93,93,93,93	0
56	MG	13	1630	1/1	0.83	0.12	-4.08	74,74,74,74	0
56	MG	1H	3414	1/1	0.98	0.15	-4.08	65,65,65,65	0
56	MG	1H	3349	1/1	0.96	0.16	-4.16	69,69,69,69	0
56	MG	1H	3057	1/1	0.96	0.19	-4.19	72,72,72,72	0
56	MG	1H	3105	1/1	0.94	0.09	-4.25	76,76,76,76	0
56	MG	1H	3112	1/1	0.94	0.14	-4.38	55,55,55,55	0
56	MG	14	3130	1/1	0.89	0.10	-4.39	71,71,71,71	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
56	MG	14	3334	1/1	0.96	0.11	-4.42	91,91,91,91	0
56	MG	14	3100	1/1	0.95	0.12	-4.43	87,87,87,87	0
56	MG	1H	3207	1/1	0.98	0.18	-4.48	60,60,60,60	0
56	MG	1H	3452	1/1	0.99	0.14	-4.51	65,65,65,65	0
56	MG	14	3218	1/1	0.92	0.13	-4.51	75,75,75,75	0
56	MG	1H	3389	1/1	0.98	0.16	-4.52	69,69,69,69	0
56	MG	13	1722	1/1	0.82	0.10	-4.55	85,85,85,85	0
56	MG	14	3351	1/1	0.98	0.12	-4.56	60,60,60,60	0
56	MG	1H	3491	1/1	0.73	0.16	-4.61	93,93,93,93	0
56	MG	1H	3206	1/1	0.77	0.14	-4.63	53,53,53,53	0
56	MG	14	3352	1/1	0.84	0.12	-4.68	74,74,74,74	0
56	MG	14	3423	1/1	0.91	0.10	-4.72	133,133,133,133	0
56	MG	1H	3358	1/1	0.91	0.15	-4.75	76,76,76,76	0
56	MG	14	3056	1/1	0.92	0.12	-4.81	92,92,92,92	0
56	MG	14	3347	1/1	0.97	0.07	-4.84	85,85,85,85	0
56	MG	1H	3347	1/1	0.95	0.13	-4.84	57,57,57,57	0
56	MG	14	3258	1/1	0.92	0.13	-4.87	104,104,104,104	0
56	MG	14	3312	1/1	0.68	0.12	-4.88	93,93,93,93	0
56	MG	1H	3136	1/1	0.88	0.15	-4.90	71,71,71,71	0
56	MG	14	3067	1/1	0.96	0.18	-4.92	73,73,73,73	0
56	MG	1H	3239	1/1	0.97	0.17	-4.93	88,88,88,88	0
56	MG	1H	3439	1/1	0.86	0.15	-4.93	92,92,92,92	0
56	MG	13	1632	1/1	0.91	0.16	-5.20	64,64,64,64	0
56	MG	14	3348	1/1	0.97	0.10	-5.20	71,71,71,71	0
56	MG	1H	3383	1/1	0.97	0.10	-5.20	67,67,67,67	0
56	MG	1H	3416	1/1	0.94	0.07	-5.25	78,78,78,78	0
56	MG	14	3240	1/1	0.94	0.10	-5.27	76,76,76,76	0
56	MG	1H	3371	1/1	0.97	0.08	-5.29	91,91,91,91	0
56	MG	1H	3404	1/1	0.94	0.09	-5.67	62,62,62,62	0
56	MG	14	3214	1/1	0.96	0.14	-5.70	70,70,70,70	0
56	MG	13	1724	1/1	0.91	0.06	-5.82	102,102,102,102	0
56	MG	1H	3411	1/1	0.99	0.12	-5.84	80,80,80,80	0
56	MG	1H	3440	1/1	0.99	0.10	-5.91	83,83,83,83	0
56	MG	1H	3352	1/1	0.98	0.14	-5.93	61,61,61,61	0
56	MG	14	3380	1/1	0.92	0.09	-5.98	97,97,97,97	0
56	MG	1H	3372	1/1	0.82	0.10	-6.28	79,79,79,79	0
56	MG	14	3105	1/1	0.97	0.14	-6.48	74,74,74,74	0
56	MG	14	3333	1/1	0.98	0.12	-6.53	73,73,73,73	0
56	MG	14	3018	1/1	0.91	0.13	-6.55	70,70,70,70	0
56	MG	1H	3443	1/1	0.99	0.17	-6.56	76,76,76,76	0
56	MG	14	3373	1/1	0.96	0.08	-6.61	81,81,81,81	0
56	MG	1H	3481	1/1	0.97	0.11	-6.74	70,70,70,70	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
56	MG	1H	3476	1/1	0.79	0.13	-7.08	93,93,93,93	0
56	MG	14	3350	1/1	0.94	0.08	-7.17	80,80,80,80	0
56	MG	14	3346	1/1	0.88	0.11	-7.40	92,92,92,92	0
56	MG	14	3412	1/1	0.85	0.08	-8.02	108,108,108,108	0
56	MG	1H	3412	1/1	0.99	0.12	-8.27	64,64,64,64	0
56	MG	14	3341	1/1	0.93	0.07	-8.39	78,78,78,78	0
56	MG	1H	3427	1/1	0.98	0.09	-8.44	75,75,75,75	0
56	MG	1H	3103	1/1	0.97	0.15	-8.99	60,60,60,60	0
56	MG	1G	1677	1/1	0.97	0.08	-9.02	91,91,91,91	0
56	MG	1H	3417	1/1	0.93	0.12	-9.21	90,90,90,90	0
56	MG	1H	3392	1/1	0.97	0.13	-9.26	65,65,65,65	0
56	MG	1H	3438	1/1	0.92	0.05	-9.42	82,82,82,82	0
56	MG	14	3335	1/1	0.99	0.12	-9.46	73,73,73,73	0
56	MG	1H	3212	1/1	0.95	0.14	-9.50	61,61,61,61	0
56	MG	1H	3384	1/1	0.95	0.07	-9.52	59,59,59,59	0
56	MG	14	3388	1/1	0.91	0.12	-10.69	85,85,85,85	0
56	MG	1H	3454	1/1	0.94	0.06	-11.11	85,85,85,85	0
56	MG	1H	3376	1/1	0.95	0.04	-14.20	96,96,96,96	0
56	MG	1H	3451	1/1	0.96	0.11	-14.82	87,87,87,87	0
56	MG	1H	3333	1/1	0.92	0.34	-	91,91,91,91	0
56	MG	13	1690	1/1	0.76	0.19	-	108,108,108,108	0
56	MG	1H	3008	1/1	0.97	0.41	-	79,79,79,79	0
56	MG	1H	3280	1/1	0.80	0.70	-	100,100,100,100	0
56	MG	1H	3144	1/1	0.88	0.23	-	80,80,80,80	0
56	MG	1H	3497	1/1	0.96	0.10	-	119,119,119,119	0
56	MG	1H	3119	1/1	0.35	0.52	-	101,101,101,101	0
56	MG	1H	3225	1/1	0.89	0.21	-	69,69,69,69	0
56	MG	14	3222	1/1	0.95	0.24	-	99,99,99,99	0
56	MG	1J	209	1/1	0.90	0.07	-	125,125,125,125	0
56	MG	1H	3450	1/1	0.96	0.12	-	66,66,66,66	0
56	MG	14	3022	1/1	0.90	0.48	-	107,107,107,107	0
56	MG	16	201	1/1	0.83	0.23	-	79,79,79,79	0
56	MG	1H	3503	1/1	0.83	0.12	-	116,116,116,116	0
56	MG	13	1644	1/1	0.90	0.12	-	96,96,96,96	0
56	MG	1H	3302	1/1	0.93	0.23	-	90,90,90,90	0
56	MG	1H	3146	1/1	0.85	0.32	-	83,83,83,83	0
56	MG	1H	3375	1/1	0.93	0.18	-	76,76,76,76	0
56	MG	14	3386	1/1	0.98	0.25	-	105,105,105,105	0
56	MG	14	3004	1/1	0.96	0.25	-	64,64,64,64	0
56	MG	1H	3227	1/1	0.97	0.26	-	123,123,123,123	0
56	MG	16	213	1/1	0.95	0.43	-	85,85,85,85	0
56	MG	1H	3140	1/1	0.82	0.52	-	101,101,101,101	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
56	MG	14	3147	1/1	0.79	0.19	-	97,97,97,97	0
56	MG	1H	3380	1/1	0.99	0.16	-	61,61,61,61	0
56	MG	13	1622	1/1	0.87	0.32	-	102,102,102,102	0
56	MG	1G	1647	1/1	0.89	0.32	-	86,86,86,86	0
56	MG	1H	3023	1/1	0.67	0.37	-	100,100,100,100	0
56	MG	13	1696	1/1	0.90	0.29	-	108,108,108,108	0
56	MG	1H	3493	1/1	0.85	0.25	-	130,130,130,130	0
56	MG	1H	3329	1/1	0.90	0.34	-	99,99,99,99	0
56	MG	1H	3095	1/1	0.95	0.43	-	74,74,74,74	0
56	MG	1G	1642	1/1	0.70	0.22	-	102,102,102,102	0
56	MG	14	3143	1/1	0.70	0.43	-	100,100,100,100	0
56	MG	1H	3357	1/1	0.98	0.07	-	57,57,57,57	0
56	MG	1H	3370	1/1	0.99	0.16	-	84,84,84,84	0
56	MG	1H	3265	1/1	0.75	0.27	-	96,96,96,96	0
56	MG	1H	3310	1/1	0.92	0.37	-	96,96,96,96	0
56	MG	13	1700	1/1	0.93	0.24	-	109,109,109,109	0
56	MG	1H	3019	1/1	0.97	0.28	-	54,54,54,54	0
56	MG	1H	3473	1/1	0.90	0.21	-	127,127,127,127	0
56	MG	13	1747	1/1	0.97	0.28	-	103,103,103,103	0
56	MG	1G	1637	1/1	0.93	0.16	-	104,104,104,104	0
56	MG	1H	3341	1/1	0.95	0.27	-	100,100,100,100	0
56	MG	14	3345	1/1	0.94	0.10	-	84,84,84,84	0
56	MG	1H	3402	1/1	0.96	0.12	-	81,81,81,81	0
56	MG	1H	3467	1/1	0.77	0.10	-	106,106,106,106	0
56	MG	13	1663	1/1	0.86	0.16	-	87,87,87,87	0
56	MG	1H	3453	1/1	0.98	0.15	-	70,70,70,70	0
56	MG	1H	3396	1/1	0.96	0.17	-	87,87,87,87	0
56	MG	1H	3367	1/1	0.95	0.13	-	82,82,82,82	0
56	MG	14	3013	1/1	0.90	0.17	-	88,88,88,88	0
56	MG	14	3319	1/1	0.71	0.29	-	118,118,118,118	0
56	MG	14	3338	1/1	0.96	0.09	-	68,68,68,68	0
56	MG	14	3313	1/1	0.75	0.19	-	124,124,124,124	0
56	MG	14	3407	1/1	0.88	0.08	-	116,116,116,116	0
56	MG	1H	3138	1/1	0.90	0.40	-	94,94,94,94	0
56	MG	14	3212	1/1	0.98	0.16	-	104,104,104,104	0
56	MG	13	1657	1/1	0.95	0.24	-	104,104,104,104	0
56	MG	14	3376	1/1	0.71	0.13	-	119,119,119,119	0
56	MG	88	202	1/1	0.87	0.30	-	77,77,77,77	0
56	MG	14	3109	1/1	0.91	0.32	-	89,89,89,89	0
56	MG	1H	3462	1/1	0.96	0.11	-	75,75,75,75	0
56	MG	14	3168	1/1	0.84	0.23	-	67,67,67,67	0
56	MG	14	3131	1/1	0.96	0.20	-	56,56,56,56	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
56	MG	13	1748	1/1	0.69	0.09	-	140,140,140,140	0
56	MG	1H	3218	1/1	0.95	0.17	-	78,78,78,78	0
56	MG	1H	3281	1/1	0.57	0.30	-	85,85,85,85	0
56	MG	E5	101	1/1	0.66	0.23	-	98,98,98,98	0
56	MG	13	1665	1/1	0.87	0.12	-	91,91,91,91	0
56	MG	14	3349	1/1	0.96	0.14	-	71,71,71,71	0
56	MG	1H	3399	1/1	0.87	0.09	-	59,59,59,59	0
56	MG	13	1719	1/1	0.86	0.13	-	95,95,95,95	0
56	MG	1H	3129	1/1	0.79	0.41	-	90,90,90,90	0
56	MG	1H	3458	1/1	0.93	0.13	-	93,93,93,93	0
56	MG	14	3401	1/1	0.94	0.06	-	120,120,120,120	0
56	MG	13	1717	1/1	0.63	0.28	-	114,114,114,114	0
56	MG	13	1736	1/1	0.99	0.17	-	102,102,102,102	0
56	MG	1H	3153	1/1	0.73	0.40	-	98,98,98,98	0
56	MG	3I	202	1/1	0.83	0.45	-	101,101,101,101	0
56	MG	1H	3488	1/1	0.85	0.15	-	114,114,114,114	0
56	MG	1H	3198	1/1	0.87	0.38	-	90,90,90,90	0
56	MG	52	201	1/1	0.83	0.12	-	107,107,107,107	0
56	MG	1H	3034	1/1	0.80	0.33	-	96,96,96,96	0
56	MG	1H	3446	1/1	0.98	0.12	-	74,74,74,74	0
56	MG	1H	3271	1/1	0.98	0.46	-	93,93,93,93	0
56	MG	1H	3386	1/1	0.94	0.16	-	82,82,82,82	0
56	MG	1H	3026	1/1	0.73	0.34	-	94,94,94,94	0
56	MG	29	302	1/1	0.56	0.35	-	108,108,108,108	0
56	MG	14	3375	1/1	0.91	0.06	-	102,102,102,102	0
56	MG	1H	3109	1/1	0.95	0.19	-	84,84,84,84	0
56	MG	13	1676	1/1	0.90	0.18	-	111,111,111,111	0
56	MG	14	3301	1/1	0.82	0.29	-	83,83,83,83	0
56	MG	1J	203	1/1	0.75	0.16	-	114,114,114,114	0
56	MG	1H	3036	1/1	0.99	0.46	-	61,61,61,61	0
56	MG	14	3421	1/1	0.94	0.05	-	130,130,130,130	0
56	MG	1H	3188	1/1	0.93	0.51	-	100,100,100,100	0
56	MG	14	3303	1/1	0.90	0.12	-	100,100,100,100	0
56	MG	14	3230	1/1	0.97	0.27	-	86,86,86,86	0
56	MG	1G	1675	1/1	0.86	0.47	-	99,99,99,99	0
56	MG	1H	3496	1/1	0.91	0.04	-	131,131,131,131	0
56	MG	1H	3339	1/1	0.86	0.33	-	90,90,90,90	0
56	MG	1H	3214	1/1	0.96	0.13	-	103,103,103,103	0
56	MG	14	3132	1/1	0.93	0.22	-	94,94,94,94	0
56	MG	14	3403	1/1	0.70	0.09	-	122,122,122,122	0
56	MG	1H	3015	1/1	0.87	0.24	-	71,71,71,71	0
56	MG	14	3207	1/1	0.91	0.14	-	68,68,68,68	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
56	MG	14	3068	1/1	0.85	0.38	-	86,86,86,86	0
56	MG	14	3383	1/1	0.90	0.16	-	98,98,98,98	0
56	MG	14	3113	1/1	0.93	0.35	-	80,80,80,80	0
56	MG	21	301	1/1	0.98	0.23	-	75,75,75,75	0
56	MG	1H	3322	1/1	0.52	0.22	-	96,96,96,96	0
56	MG	16	210	1/1	0.95	0.45	-	107,107,107,107	0
56	MG	1H	3210	1/1	0.89	0.17	-	69,69,69,69	0
56	MG	1H	3286	1/1	0.84	0.26	-	85,85,85,85	0
56	MG	1G	1663	1/1	0.83	0.20	-	92,92,92,92	0
56	MG	13	1677	1/1	0.91	0.39	-	98,98,98,98	0
56	MG	1H	3466	1/1	0.90	0.12	-	94,94,94,94	0
56	MG	14	3031	1/1	0.96	0.29	-	76,76,76,76	0
56	MG	1H	3290	1/1	0.86	0.49	-	92,92,92,92	0
56	MG	2K	102	1/1	0.94	0.23	-	97,97,97,97	0
56	MG	1H	3083	1/1	0.87	0.32	-	92,92,92,92	0
56	MG	15	201	1/1	0.84	0.54	-	102,102,102,102	0
56	MG	1H	3419	1/1	0.95	0.16	-	94,94,94,94	0
56	MG	14	3114	1/1	0.95	0.39	-	97,97,97,97	0
56	MG	5E	201	1/1	0.87	0.20	-	104,104,104,104	0
56	MG	1G	1611	1/1	0.92	0.18	-	105,105,105,105	0
56	MG	1H	3262	1/1	0.81	0.30	-	89,89,89,89	0
56	MG	1G	1643	1/1	0.74	0.30	-	104,104,104,104	0
56	MG	14	3287	1/1	0.90	0.20	-	95,95,95,95	0
56	MG	1G	1613	1/1	0.92	0.22	-	117,117,117,117	0
56	MG	25	202	1/1	0.87	0.33	-	125,125,125,125	0
56	MG	14	3190	1/1	0.94	0.22	-	124,124,124,124	0
56	MG	1H	3060	1/1	0.98	0.23	-	74,74,74,74	0
56	MG	14	3224	1/1	0.93	0.09	-	80,80,80,80	0
56	MG	14	3203	1/1	0.70	0.29	-	81,81,81,81	0
56	MG	14	3389	1/1	0.96	0.17	-	72,72,72,72	0
56	MG	14	3173	1/1	0.82	0.28	-	87,87,87,87	0
56	MG	14	3422	1/1	0.79	0.09	-	132,132,132,132	0
56	MG	1H	3309	1/1	0.61	0.45	-	96,96,96,96	0
56	MG	13	1750	1/1	0.45	0.35	-	131,131,131,131	0
56	MG	14	3124	1/1	0.97	0.26	-	85,85,85,85	0
56	MG	1H	3469	1/1	0.90	0.07	-	106,106,106,106	0
56	MG	1H	3338	1/1	0.92	0.27	-	110,110,110,110	0
56	MG	14	3149	1/1	0.89	0.21	-	79,79,79,79	0
56	MG	1G	1668	1/1	0.83	0.22	-	111,111,111,111	0
56	MG	1G	1616	1/1	0.95	0.18	-	149,149,149,149	0
56	MG	1G	1672	1/1	0.95	0.26	-	126,126,126,126	0
56	MG	14	3039	1/1	0.92	0.26	-	68,68,68,68	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
56	MG	1G	1612	1/1	0.95	0.15	-	110,110,110,110	0
56	MG	13	1681	1/1	0.96	0.25	-	107,107,107,107	0
56	MG	1H	3293	1/1	0.93	0.17	-	85,85,85,85	0
56	MG	13	1606	1/1	0.60	0.33	-	115,115,115,115	0
56	MG	45	202	1/1	0.88	0.53	-	82,82,82,82	0
56	MG	14	3026	1/1	0.68	0.39	-	92,92,92,92	0
56	MG	1H	3195	1/1	0.97	0.11	-	94,94,94,94	0
56	MG	1H	3061	1/1	0.85	0.34	-	85,85,85,85	0
56	MG	1H	3038	1/1	0.96	0.26	-	58,58,58,58	0
56	MG	13	1645	1/1	0.93	0.38	-	86,86,86,86	0
56	MG	16	202	1/1	0.94	0.28	-	73,73,73,73	0
56	MG	1G	1652	1/1	0.88	0.26	-	103,103,103,103	0
56	MG	13	1634	1/1	0.98	0.30	-	85,85,85,85	0
56	MG	1H	3185	1/1	0.84	0.38	-	88,88,88,88	0
56	MG	1G	1610	1/1	0.91	0.13	-	115,115,115,115	0
56	MG	14	3324	1/1	0.82	0.40	-	96,96,96,96	0
56	MG	1H	3179	1/1	0.72	0.40	-	96,96,96,96	0
56	MG	1G	1632	1/1	0.77	0.41	-	116,116,116,116	0
56	MG	14	3201	1/1	0.63	0.30	-	111,111,111,111	0
56	MG	14	3321	1/1	0.92	0.26	-	99,99,99,99	0
56	MG	1H	3442	1/1	0.93	0.12	-	57,57,57,57	0
56	MG	14	3161	1/1	0.90	0.19	-	104,104,104,104	0
56	MG	14	3193	1/1	0.83	0.09	-	89,89,89,89	0
56	MG	1H	3435	1/1	0.99	0.15	-	72,72,72,72	0
56	MG	14	3126	1/1	0.95	0.29	-	103,103,103,103	0
56	MG	1H	3492	1/1	0.86	0.23	-	120,120,120,120	0
56	MG	1H	3226	1/1	0.97	0.26	-	88,88,88,88	0
56	MG	14	3366	1/1	0.99	0.15	-	88,88,88,88	0
56	MG	1G	1683	1/1	0.96	0.14	-	129,129,129,129	0
56	MG	1H	3336	1/1	0.93	0.30	-	102,102,102,102	0
56	MG	78	201	1/1	0.92	0.15	-	67,67,67,67	0
56	MG	14	3417	1/1	0.93	0.06	-	148,148,148,148	0
56	MG	13	1740	1/1	0.85	0.17	-	106,106,106,106	0
56	MG	1H	3337	1/1	0.65	0.42	-	106,106,106,106	0
56	MG	11	302	1/1	0.90	0.25	-	66,66,66,66	0
56	MG	1H	3264	1/1	0.93	0.29	-	96,96,96,96	0
56	MG	1H	3426	1/1	0.91	0.13	-	92,92,92,92	0
56	MG	1H	3395	1/1	0.65	0.17	-	118,118,118,118	0
56	MG	1H	3035	1/1	0.93	0.20	-	55,55,55,55	0
56	MG	14	3034	1/1	0.98	0.23	-	69,69,69,69	0
56	MG	1G	1676	1/1	0.95	0.35	-	121,121,121,121	0
56	MG	14	3191	1/1	0.97	0.21	-	92,92,92,92	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
56	MG	14	3316	1/1	0.87	0.54	-	103,103,103,103	0
56	MG	14	3108	1/1	0.96	0.15	-	92,92,92,92	0
56	MG	1H	3016	1/1	0.98	0.31	-	66,66,66,66	0
56	MG	1H	3482	1/1	0.94	0.10	-	119,119,119,119	0
56	MG	1H	3031	1/1	0.89	0.33	-	89,89,89,89	0
56	MG	1H	3263	1/1	0.86	0.22	-	106,106,106,106	0
56	MG	1H	3247	1/1	0.93	0.28	-	81,81,81,81	0
56	MG	1H	3236	1/1	0.95	0.26	-	77,77,77,77	0
56	MG	32	301	1/1	0.76	0.46	-	120,120,120,120	0
56	MG	1H	3191	1/1	0.91	0.34	-	87,87,87,87	0
56	MG	1H	3348	1/1	0.97	0.14	-	84,84,84,84	0
56	MG	14	3196	1/1	0.77	0.34	-	99,99,99,99	0
56	MG	14	3250	1/1	0.89	0.23	-	86,86,86,86	0
56	MG	1H	3093	1/1	0.96	0.16	-	81,81,81,81	0
56	MG	1G	1631	1/1	0.95	0.23	-	113,113,113,113	0
56	MG	13	1669	1/1	0.70	0.27	-	110,110,110,110	0
56	MG	14	3286	1/1	0.69	0.15	-	92,92,92,92	0
56	MG	1H	3480	1/1	0.87	0.14	-	109,109,109,109	0
56	MG	G8	201	1/1	0.95	0.11	-	83,83,83,83	0
56	MG	6A	101	1/1	0.84	0.21	-	138,138,138,138	0
56	MG	1H	3122	1/1	0.84	0.30	-	94,94,94,94	0
56	MG	14	3177	1/1	0.89	0.32	-	111,111,111,111	0
56	MG	1H	3422	1/1	0.94	0.15	-	91,91,91,91	0
56	MG	1H	3343	1/1	0.68	0.36	-	92,92,92,92	0
56	MG	13	1702	1/1	0.79	0.23	-	109,109,109,109	0
56	MG	1H	3199	1/1	0.84	0.51	-	99,99,99,99	0
56	MG	1H	3253	1/1	0.97	0.19	-	91,91,91,91	0
56	MG	13	1636	1/1	0.62	0.30	-	100,100,100,100	0
56	MG	1G	1658	1/1	0.95	0.21	-	110,110,110,110	0
56	MG	14	3367	1/1	0.92	0.06	-	111,111,111,111	0
56	MG	14	3135	1/1	0.93	0.31	-	101,101,101,101	0
56	MG	14	3400	1/1	0.93	0.22	-	95,95,95,95	0
56	MG	13	1624	1/1	0.87	0.22	-	95,95,95,95	0
56	MG	13	1673	1/1	0.95	0.28	-	91,91,91,91	0
56	MG	1H	3400	1/1	0.94	0.10	-	61,61,61,61	0
56	MG	13	1734	1/1	0.81	0.15	-	133,133,133,133	0
56	MG	1H	3312	1/1	0.91	0.37	-	79,79,79,79	0
56	MG	1H	3413	1/1	0.97	0.15	-	77,77,77,77	0
56	MG	1H	3167	1/1	0.93	0.15	-	92,92,92,92	0
56	MG	1G	1646	1/1	0.96	0.15	-	102,102,102,102	0
56	MG	1G	1606	1/1	0.90	0.27	-	129,129,129,129	0
56	MG	13	1744	1/1	0.68	0.41	-	132,132,132,132	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
56	MG	13	1739	1/1	0.47	0.20	-	120,120,120,120	0
56	MG	14	3281	1/1	0.75	0.55	-	105,105,105,105	0
56	MG	1H	3460	1/1	0.94	0.11	-	110,110,110,110	0
56	MG	1H	3455	1/1	0.94	0.22	-	99,99,99,99	0
56	MG	14	3093	1/1	0.85	0.51	-	103,103,103,103	0
56	MG	14	3251	1/1	0.71	0.10	-	105,105,105,105	0
56	MG	14	3320	1/1	0.79	0.22	-	96,96,96,96	0
56	MG	14	3054	1/1	0.96	0.12	-	93,93,93,93	0
56	MG	14	3241	1/1	0.94	0.31	-	92,92,92,92	0
56	MG	14	3339	1/1	0.99	0.12	-	74,74,74,74	0
56	MG	13	1714	1/1	0.70	0.22	-	101,101,101,101	0
56	MG	1H	3295	1/1	0.88	0.50	-	92,92,92,92	0
56	MG	1H	3217	1/1	0.93	0.12	-	91,91,91,91	0
56	MG	1H	3118	1/1	0.88	0.29	-	76,76,76,76	0
56	MG	14	3402	1/1	0.96	0.11	-	99,99,99,99	0
56	MG	13	1687	1/1	0.84	0.28	-	109,109,109,109	0
56	MG	1H	3465	1/1	0.98	0.10	-	83,83,83,83	0
56	MG	14	3134	1/1	0.97	0.10	-	103,103,103,103	0
56	MG	14	3166	1/1	0.89	0.14	-	106,106,106,106	0
56	MG	13	1745	1/1	0.61	0.39	-	121,121,121,121	0
56	MG	14	3006	1/1	0.94	0.21	-	99,99,99,99	0
56	MG	1G	1686	1/1	0.82	0.07	-	129,129,129,129	0
56	MG	14	3200	1/1	0.85	0.18	-	85,85,85,85	0
56	MG	14	3099	1/1	0.95	0.14	-	106,106,106,106	0
56	MG	1H	3090	1/1	0.95	0.46	-	91,91,91,91	0
56	MG	14	3360	1/1	0.91	0.06	-	92,92,92,92	0
56	MG	14	3164	1/1	0.88	0.21	-	91,91,91,91	0
56	MG	14	3298	1/1	0.95	0.42	-	115,115,115,115	0
56	MG	14	3085	1/1	0.97	0.18	-	81,81,81,81	0
56	MG	14	3282	1/1	0.88	0.31	-	90,90,90,90	0
56	MG	1H	3423	1/1	0.84	0.14	-	102,102,102,102	0
56	MG	1H	3424	1/1	0.94	0.09	-	119,119,119,119	0
56	MG	14	3025	1/1	0.62	0.17	-	116,116,116,116	0
56	MG	14	3405	1/1	0.81	0.22	-	102,102,102,102	0
56	MG	14	3088	1/1	0.96	0.23	-	66,66,66,66	0
56	MG	13	1720	1/1	0.43	0.24	-	104,104,104,104	0
56	MG	1H	3365	1/1	0.98	0.14	-	69,69,69,69	0
56	MG	14	3012	1/1	0.92	0.23	-	94,94,94,94	0
56	MG	1H	3079	1/1	0.99	0.33	-	83,83,83,83	0
56	MG	1H	3373	1/1	0.98	0.12	-	63,63,63,63	0
56	MG	1H	3278	1/1	0.92	0.29	-	101,101,101,101	0
56	MG	1H	3072	1/1	0.79	0.29	-	80,80,80,80	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
56	MG	1H	3468	1/1	0.97	0.07	-	95,95,95,95	0
56	MG	14	3075	1/1	0.98	0.23	-	81,81,81,81	0
56	MG	1H	3059	1/1	0.98	0.49	-	92,92,92,92	0
56	MG	1H	3028	1/1	0.49	0.43	-	115,115,115,115	0
56	MG	1H	3483	1/1	0.94	0.07	-	110,110,110,110	0
56	MG	14	3331	1/1	0.99	0.29	-	112,112,112,112	0
56	MG	1J	206	1/1	0.82	0.12	-	123,123,123,123	0
56	MG	1H	3049	1/1	0.93	0.33	-	60,60,60,60	0
56	MG	13	1715	1/1	0.94	0.30	-	108,108,108,108	0
56	MG	1H	3161	1/1	0.94	0.36	-	82,82,82,82	0
56	MG	1H	3405	1/1	0.95	0.13	-	86,86,86,86	0
56	MG	14	3063	1/1	0.93	0.18	-	104,104,104,104	0
56	MG	1H	3335	1/1	0.87	0.42	-	85,85,85,85	0
56	MG	1H	3409	1/1	0.99	0.18	-	76,76,76,76	0
56	MG	1H	3246	1/1	0.84	0.22	-	78,78,78,78	0
56	MG	14	3101	1/1	0.98	0.15	-	80,80,80,80	0
56	MG	1G	1617	1/1	0.92	0.27	-	110,110,110,110	0
56	MG	14	3354	1/1	0.88	0.06	-	126,126,126,126	0
56	MG	14	3083	1/1	0.92	0.26	-	81,81,81,81	0
56	MG	14	3234	1/1	0.94	0.55	-	98,98,98,98	0
56	MG	1H	3354	1/1	0.92	0.14	-	64,64,64,64	0
56	MG	14	3152	1/1	0.86	0.23	-	80,80,80,80	0
56	MG	1H	3238	1/1	0.90	0.32	-	84,84,84,84	0
56	MG	1H	3111	1/1	0.83	0.43	-	91,91,91,91	0
56	MG	14	3249	1/1	0.87	0.27	-	97,97,97,97	0
56	MG	P8	101	1/1	0.85	0.41	-	84,84,84,84	0
56	MG	14	3288	1/1	0.56	0.53	-	107,107,107,107	0
56	MG	1H	3092	1/1	0.85	0.30	-	70,70,70,70	0
56	MG	14	3133	1/1	0.68	0.13	-	97,97,97,97	0
56	MG	1H	3018	1/1	0.97	0.40	-	79,79,79,79	0
56	MG	14	3326	1/1	0.81	0.39	-	117,117,117,117	0
56	MG	1H	3317	1/1	0.86	0.79	-	85,85,85,85	0
56	MG	13	1633	1/1	0.98	0.15	-	75,75,75,75	0
56	MG	1H	3142	1/1	0.41	0.41	-	88,88,88,88	0
56	MG	1H	3187	1/1	0.96	0.24	-	89,89,89,89	0
56	MG	14	3145	1/1	0.97	0.25	-	105,105,105,105	0
56	MG	1H	3231	1/1	0.93	0.49	-	90,90,90,90	0
56	MG	1H	3283	1/1	0.84	0.38	-	112,112,112,112	0
56	MG	1H	3106	1/1	0.76	0.34	-	95,95,95,95	0
56	MG	1G	1649	1/1	0.73	0.27	-	105,105,105,105	0
56	MG	14	3175	1/1	0.83	0.38	-	99,99,99,99	0
56	MG	1H	3192	1/1	0.96	0.11	-	78,78,78,78	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
56	MG	14	3330	1/1	0.83	0.20	-	105,105,105,105	0
56	MG	14	3046	1/1	0.97	0.24	-	88,88,88,88	0
56	MG	13	1615	1/1	0.94	0.24	-	100,100,100,100	0
56	MG	1H	3022	1/1	0.99	0.38	-	64,64,64,64	0
56	MG	13	1688	1/1	0.72	0.33	-	93,93,93,93	0
56	MG	14	3297	1/1	0.72	0.24	-	96,96,96,96	0
56	MG	1H	3391	1/1	0.84	0.09	-	96,96,96,96	0
56	MG	14	3112	1/1	0.90	0.12	-	98,98,98,98	0
56	MG	14	3424	1/1	0.91	0.43	-	116,116,116,116	0
56	MG	1H	3313	1/1	0.57	0.24	-	93,93,93,93	0
56	MG	1H	3100	1/1	0.91	0.27	-	86,86,86,86	0
56	MG	2K	103	1/1	0.95	0.10	-	98,98,98,98	0
56	MG	1H	3050	1/1	0.93	0.27	-	80,80,80,80	0
56	MG	1H	3308	1/1	0.90	0.25	-	90,90,90,90	0
56	MG	1H	3306	1/1	0.68	0.61	-	81,81,81,81	0
56	MG	1H	3318	1/1	0.84	0.30	-	87,87,87,87	0
56	MG	1H	3223	1/1	0.87	0.18	-	82,82,82,82	0
56	MG	14	3129	1/1	0.96	0.20	-	78,78,78,78	0
56	MG	1H	3345	1/1	0.89	0.31	-	100,100,100,100	0
56	MG	14	3309	1/1	0.92	0.49	-	87,87,87,87	0
56	MG	14	3066	1/1	0.87	0.30	-	96,96,96,96	0
56	MG	14	3391	1/1	0.46	0.10	-	124,124,124,124	0
56	MG	1H	3200	1/1	0.57	0.35	-	104,104,104,104	0
56	MG	14	3187	1/1	0.88	0.36	-	100,100,100,100	0
56	MG	14	3117	1/1	0.80	0.45	-	98,98,98,98	0
56	MG	1G	1654	1/1	0.94	0.38	-	94,94,94,94	0
56	MG	1H	3194	1/1	0.92	0.20	-	104,104,104,104	0
56	MG	14	3342	1/1	0.99	0.12	-	80,80,80,80	0
56	MG	1H	3284	1/1	0.87	0.26	-	70,70,70,70	0
56	MG	1H	3296	1/1	0.75	0.21	-	86,86,86,86	0
56	MG	1H	3221	1/1	0.94	0.38	-	99,99,99,99	0
56	MG	14	3414	1/1	0.98	0.06	-	93,93,93,93	0
56	MG	14	3087	1/1	0.96	0.16	-	82,82,82,82	0
56	MG	14	3263	1/1	0.93	0.25	-	88,88,88,88	0
56	MG	1H	3208	1/1	0.94	0.09	-	63,63,63,63	0
56	MG	14	3104	1/1	0.84	0.27	-	81,81,81,81	0
56	MG	1H	3268	1/1	0.90	0.50	-	125,125,125,125	0
56	MG	14	3125	1/1	0.89	0.27	-	101,101,101,101	0
56	MG	14	3148	1/1	0.86	0.11	-	117,117,117,117	0
56	MG	1G	1670	1/1	0.64	0.40	-	138,138,138,138	0
56	MG	13	1698	1/1	0.94	0.35	-	95,95,95,95	0
56	MG	1G	1669	1/1	0.84	0.28	-	112,112,112,112	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
56	MG	14	3239	1/1	0.90	0.23	-	166,166,166,166	0
56	MG	1H	3196	1/1	0.95	0.70	-	110,110,110,110	0
56	MG	1H	3276	1/1	0.65	0.44	-	99,99,99,99	0
56	MG	13	1716	1/1	0.68	0.18	-	111,111,111,111	0
56	MG	1H	3080	1/1	0.94	0.31	-	86,86,86,86	0
56	MG	14	3275	1/1	0.79	0.28	-	105,105,105,105	0
56	MG	1H	3157	1/1	0.71	0.51	-	104,104,104,104	0
56	MG	13	1626	1/1	0.94	0.20	-	109,109,109,109	0
56	MG	1H	3330	1/1	0.69	0.56	-	104,104,104,104	0
56	MG	14	3092	1/1	0.94	0.16	-	109,109,109,109	0
56	MG	1H	3323	1/1	0.87	0.11	-	110,110,110,110	0
56	MG	1H	3485	1/1	0.82	0.08	-	121,121,121,121	0
56	MG	1H	3012	1/1	0.84	0.50	-	93,93,93,93	0
56	MG	1H	3039	1/1	0.97	0.39	-	55,55,55,55	0
56	MG	14	3355	1/1	0.90	0.18	-	102,102,102,102	0
56	MG	1H	3058	1/1	0.94	0.20	-	107,107,107,107	0
56	MG	1G	1665	1/1	0.87	0.30	-	110,110,110,110	0
56	MG	1H	3499	1/1	0.81	0.11	-	110,110,110,110	0
56	MG	14	3136	1/1	0.84	0.10	-	112,112,112,112	0
56	MG	14	3017	1/1	0.90	0.15	-	91,91,91,91	0
56	MG	1H	3032	1/1	0.93	0.25	-	78,78,78,78	0
56	MG	14	3396	1/1	0.91	0.16	-	119,119,119,119	0
56	MG	1G	1619	1/1	0.96	0.21	-	72,72,72,72	0
56	MG	1H	3299	1/1	0.97	0.17	-	71,71,71,71	0
56	MG	14	3183	1/1	0.51	0.32	-	110,110,110,110	0
56	MG	14	3178	1/1	0.78	0.44	-	95,95,95,95	0
56	MG	14	3397	1/1	0.97	0.13	-	101,101,101,101	0
56	MG	14	3181	1/1	0.61	0.21	-	100,100,100,100	0
56	MG	14	3019	1/1	0.66	0.22	-	96,96,96,96	0
56	MG	14	3285	1/1	0.75	0.30	-	91,91,91,91	0
56	MG	1H	3436	1/1	0.96	0.10	-	79,79,79,79	0
56	MG	1H	3007	1/1	0.86	0.33	-	97,97,97,97	0
56	MG	1G	1684	1/1	0.84	0.12	-	124,124,124,124	0
56	MG	1H	3229	1/1	0.96	0.23	-	83,83,83,83	0
56	MG	13	1749	1/1	0.70	0.23	-	121,121,121,121	0
56	MG	31	301	1/1	0.86	0.10	-	92,92,92,92	0
56	MG	13	1661	1/1	0.93	0.29	-	91,91,91,91	0
56	MG	1H	3494	1/1	0.87	0.36	-	121,121,121,121	0
56	MG	14	3226	1/1	0.96	0.21	-	123,123,123,123	0
56	MG	1H	3070	1/1	0.98	0.54	-	125,125,125,125	0
56	MG	1J	208	1/1	0.81	0.11	-	121,121,121,121	0
56	MG	1H	3319	1/1	0.93	0.27	-	90,90,90,90	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
56	MG	14	3188	1/1	0.98	0.31	-	91,91,91,91	0
56	MG	14	3361	1/1	0.97	0.17	-	86,86,86,86	0
56	MG	13	1727	1/1	0.89	0.24	-	121,121,121,121	0
56	MG	14	3050	1/1	0.96	0.13	-	60,60,60,60	0
56	MG	1G	1604	1/1	0.77	0.23	-	142,142,142,142	0
56	MG	1H	3429	1/1	0.94	0.08	-	102,102,102,102	0
56	MG	14	3023	1/1	0.88	0.15	-	86,86,86,86	0
56	MG	14	3255	1/1	0.77	0.41	-	99,99,99,99	0
56	MG	1H	3374	1/1	0.93	0.16	-	71,71,71,71	0
56	MG	14	3036	1/1	0.97	0.15	-	67,67,67,67	0
56	MG	13	1699	1/1	0.84	0.50	-	102,102,102,102	0
56	MG	14	3043	1/1	0.98	0.21	-	101,101,101,101	0
56	MG	14	3317	1/1	0.73	0.28	-	98,98,98,98	0
56	MG	1H	3303	1/1	0.93	0.30	-	76,76,76,76	0
56	MG	14	3404	1/1	0.96	0.08	-	97,97,97,97	0
56	MG	1H	3377	1/1	0.96	0.17	-	97,97,97,97	0
56	MG	1H	3062	1/1	0.74	0.23	-	101,101,101,101	0
56	MG	1H	3021	1/1	0.95	0.32	-	79,79,79,79	0
56	MG	14	3280	1/1	0.81	0.31	-	95,95,95,95	0
56	MG	1H	3506	1/1	0.96	0.33	-	57,57,57,57	0
56	MG	14	3315	1/1	0.75	0.12	-	92,92,92,92	0
56	MG	14	3356	1/1	0.71	0.25	-	118,118,118,118	0
56	MG	1H	3224	1/1	0.98	0.28	-	82,82,82,82	0
56	MG	1H	3502	1/1	0.89	0.29	-	118,118,118,118	0
56	MG	1H	3030	1/1	0.86	0.19	-	79,79,79,79	0
56	MG	14	3294	1/1	0.72	0.14	-	91,91,91,91	0
56	MG	1H	3201	1/1	0.93	0.32	-	121,121,121,121	0
56	MG	14	3343	1/1	0.88	0.12	-	89,89,89,89	0
56	MG	1H	3137	1/1	0.73	0.53	-	87,87,87,87	0
56	MG	1H	3344	1/1	0.85	0.42	-	95,95,95,95	0
56	MG	1H	3505	1/1	0.71	0.15	-	112,112,112,112	0
56	MG	1H	3325	1/1	0.94	0.13	-	86,86,86,86	0
56	MG	14	3035	1/1	0.99	0.27	-	79,79,79,79	0
56	MG	13	1672	1/1	0.88	0.20	-	100,100,100,100	0
56	MG	13	1627	1/1	0.66	0.28	-	101,101,101,101	0
56	MG	1H	3478	1/1	0.87	0.09	-	118,118,118,118	0
56	MG	14	3086	1/1	0.96	0.28	-	84,84,84,84	0
56	MG	14	3028	1/1	0.90	0.14	-	98,98,98,98	0
56	MG	14	3003	1/1	0.94	0.13	-	73,73,73,73	0
56	MG	1H	3074	1/1	0.80	0.41	-	80,80,80,80	0
56	MG	14	3378	1/1	0.96	0.16	-	101,101,101,101	0
56	MG	1H	3420	1/1	0.92	0.10	-	84,84,84,84	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
56	MG	1G	1650	1/1	0.86	0.36	-	101,101,101,101	0
56	MG	1G	1638	1/1	0.84	0.42	-	95,95,95,95	0
56	MG	14	3158	1/1	0.77	0.15	-	93,93,93,93	0
56	MG	1H	3418	1/1	0.93	0.06	-	134,134,134,134	0
56	MG	1H	3433	1/1	0.95	0.09	-	76,76,76,76	0
56	MG	1H	3394	1/1	0.97	0.09	-	62,62,62,62	0
56	MG	13	1646	1/1	0.90	0.24	-	104,104,104,104	0
56	MG	14	3408	1/1	0.98	0.08	-	79,79,79,79	0
56	MG	1H	3407	1/1	0.77	0.07	-	115,115,115,115	0
56	MG	14	3377	1/1	0.98	0.08	-	82,82,82,82	0
56	MG	14	3029	1/1	0.69	0.22	-	97,97,97,97	0
56	MG	14	3030	1/1	0.74	0.22	-	115,115,115,115	0
56	MG	14	3283	1/1	0.87	0.21	-	98,98,98,98	0
56	MG	14	3077	1/1	0.99	0.28	-	80,80,80,80	0
56	MG	I8	102	1/1	0.90	0.41	-	64,64,64,64	0
56	MG	78	202	1/1	0.93	0.20	-	97,97,97,97	0
56	MG	13	1693	1/1	0.78	0.21	-	103,103,103,103	0
56	MG	13	1670	1/1	0.86	0.33	-	88,88,88,88	0
56	MG	13	1629	1/1	0.62	0.45	-	106,106,106,106	0
56	MG	1H	3428	1/1	0.75	0.06	-	123,123,123,123	0
56	MG	14	3399	1/1	0.94	0.08	-	78,78,78,78	0
56	MG	1H	3183	1/1	0.93	0.29	-	83,83,83,83	0
56	MG	1H	3274	1/1	0.90	0.57	-	100,100,100,100	0
56	MG	14	3127	1/1	0.94	0.06	-	87,87,87,87	0
56	MG	14	3151	1/1	0.83	0.24	-	91,91,91,91	0
56	MG	1H	3495	1/1	0.92	0.32	-	91,91,91,91	0
56	MG	1H	3501	1/1	0.92	0.17	-	107,107,107,107	0
56	MG	1H	3363	1/1	0.96	0.21	-	71,71,71,71	0
56	MG	1H	3346	1/1	0.75	0.61	-	106,106,106,106	0
56	MG	1H	3149	1/1	0.92	0.54	-	99,99,99,99	0
56	MG	1H	3182	1/1	0.85	0.36	-	74,74,74,74	0
56	MG	14	3024	1/1	0.93	0.28	-	73,73,73,73	0
56	MG	1G	1629	1/1	0.55	0.18	-	124,124,124,124	0
56	MG	1H	3381	1/1	0.88	0.09	-	89,89,89,89	0
56	MG	14	3096	1/1	0.99	0.15	-	77,77,77,77	0
56	MG	14	3140	1/1	0.92	0.34	-	93,93,93,93	0
56	MG	13	1683	1/1	0.57	0.28	-	106,106,106,106	0
56	MG	1G	1605	1/1	0.90	0.17	-	98,98,98,98	0
56	MG	14	3277	1/1	0.88	0.20	-	90,90,90,90	0
56	MG	13	1658	1/1	0.57	0.44	-	117,117,117,117	0
56	MG	1H	3397	1/1	0.97	0.15	-	77,77,77,77	0
56	MG	1H	3415	1/1	0.75	0.07	-	99,99,99,99	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
56	MG	13	1730	1/1	0.95	0.03	-	113,113,113,113	0
56	MG	14	3302	1/1	0.91	0.19	-	92,92,92,92	0
56	MG	1H	3279	1/1	0.90	0.25	-	80,80,80,80	0
56	MG	13	1616	1/1	0.95	0.10	-	82,82,82,82	0
56	MG	14	3311	1/1	0.76	0.17	-	89,89,89,89	0
56	MG	1G	1648	1/1	0.83	0.23	-	100,100,100,100	0
56	MG	14	3273	1/1	0.88	0.14	-	87,87,87,87	0
56	MG	14	3210	1/1	0.91	0.37	-	98,98,98,98	0
56	MG	13	1610	1/1	0.97	0.20	-	79,79,79,79	0
56	MG	1H	3065	1/1	0.96	0.27	-	71,71,71,71	0
56	MG	14	3144	1/1	0.81	0.35	-	93,93,93,93	0
56	MG	1H	3291	1/1	0.92	0.29	-	77,77,77,77	0
56	MG	1H	3010	1/1	0.95	0.34	-	83,83,83,83	0
56	MG	1G	1602	1/1	0.85	0.38	-	118,118,118,118	0
56	MG	14	3291	1/1	0.71	0.18	-	115,115,115,115	0
56	MG	1H	3398	1/1	0.94	0.09	-	92,92,92,92	0
56	MG	1H	3305	1/1	0.80	0.61	-	104,104,104,104	0
56	MG	1G	1666	1/1	0.90	0.36	-	89,89,89,89	0
56	MG	1H	3406	1/1	0.98	0.16	-	77,77,77,77	0
56	MG	25	201	1/1	0.83	0.43	-	111,111,111,111	0
56	MG	C5	201	1/1	0.86	0.23	-	123,123,123,123	0
56	MG	13	1691	1/1	0.87	0.30	-	95,95,95,95	0
56	MG	14	3089	1/1	0.97	0.21	-	85,85,85,85	0
56	MG	13	1692	1/1	0.56	0.22	-	102,102,102,102	0
56	MG	14	3305	1/1	0.87	0.14	-	79,79,79,79	0
56	MG	1H	3047	1/1	0.95	0.30	-	65,65,65,65	0
56	MG	1H	3474	1/1	0.94	0.20	-	124,124,124,124	0
56	MG	14	3374	1/1	0.83	0.08	-	113,113,113,113	0
56	MG	14	3027	1/1	0.93	0.20	-	84,84,84,84	0
56	MG	14	3159	1/1	0.91	0.13	-	112,112,112,112	0
56	MG	1H	3202	1/1	0.78	0.37	-	93,93,93,93	0
56	MG	1H	3175	1/1	0.77	0.35	-	93,93,93,93	0
56	MG	14	3276	1/1	0.74	1.34	-	103,103,103,103	0
56	MG	14	3179	1/1	0.95	0.31	-	96,96,96,96	0
56	MG	1H	3487	1/1	0.83	0.29	-	126,126,126,126	0
56	MG	1H	3174	1/1	0.85	0.31	-	78,78,78,78	0
56	MG	14	3185	1/1	0.84	0.20	-	90,90,90,90	0
56	MG	1J	210	1/1	0.94	0.06	-	109,109,109,109	0
56	MG	13	1705	1/1	0.98	0.14	-	98,98,98,98	0
56	MG	1H	3073	1/1	0.94	0.24	-	88,88,88,88	0
56	MG	13	1653	1/1	0.95	0.21	-	78,78,78,78	0
56	MG	1H	3334	1/1	0.91	0.50	-	99,99,99,99	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
56	MG	14	3123	1/1	0.87	0.16	-	75,75,75,75	0
56	MG	14	3037	1/1	0.90	0.20	-	79,79,79,79	0
56	MG	14	3213	1/1	0.93	0.17	-	113,113,113,113	0
56	MG	13	1746	1/1	0.97	0.07	-	110,110,110,110	0
56	MG	1H	3267	1/1	0.79	0.34	-	84,84,84,84	0
56	MG	1H	3311	1/1	0.68	0.48	-	109,109,109,109	0
56	MG	1H	3282	1/1	0.88	0.14	-	107,107,107,107	0
56	MG	1J	202	1/1	0.86	0.12	-	157,157,157,157	0
56	MG	14	3410	1/1	0.94	0.30	-	96,96,96,96	0
56	MG	1H	3134	1/1	0.96	0.13	-	57,57,57,57	0
56	MG	14	3059	1/1	0.94	0.27	-	74,74,74,74	0
56	MG	14	3295	1/1	0.61	0.62	-	95,95,95,95	0
56	MG	13	1631	1/1	0.51	0.70	-	102,102,102,102	0
56	MG	1H	3086	1/1	0.87	0.39	-	72,72,72,72	0
56	MG	1G	1661	1/1	0.90	0.26	-	105,105,105,105	0
56	MG	2K	104	1/1	0.87	0.44	-	90,90,90,90	0
56	MG	1H	3484	1/1	0.98	0.09	-	79,79,79,79	0
56	MG	1G	1615	1/1	0.86	0.23	-	115,115,115,115	0
56	MG	1H	3126	1/1	0.82	0.32	-	79,79,79,79	0
56	MG	14	3174	1/1	0.88	0.18	-	93,93,93,93	0
56	MG	1H	3135	1/1	0.70	0.26	-	111,111,111,111	0
56	MG	1H	3158	1/1	0.62	0.46	-	104,104,104,104	0
56	MG	14	3292	1/1	0.93	0.17	-	90,90,90,90	0
56	MG	13	1605	1/1	0.86	0.18	-	93,93,93,93	0
56	MG	14	3225	1/1	0.94	0.20	-	65,65,65,65	0
56	MG	14	3325	1/1	0.83	0.34	-	89,89,89,89	0
56	MG	14	3167	1/1	0.79	0.27	-	110,110,110,110	0
56	MG	14	3245	1/1	0.96	0.18	-	71,71,71,71	0
56	MG	13	1656	1/1	0.74	0.42	-	103,103,103,103	0
56	MG	14	3359	1/1	0.95	0.13	-	114,114,114,114	0
56	MG	1H	3321	1/1	0.85	0.27	-	99,99,99,99	0
56	MG	14	3384	1/1	0.95	0.14	-	60,60,60,60	0
56	MG	14	3020	1/1	0.54	0.51	-	92,92,92,92	0
56	MG	14	3372	1/1	0.70	0.12	-	134,134,134,134	0
56	MG	1H	3316	1/1	0.90	0.57	-	94,94,94,94	0
56	MG	1H	3498	1/1	0.39	0.20	-	119,119,119,119	0
56	MG	M5	101	1/1	0.56	0.25	-	94,94,94,94	0
56	MG	1H	3143	1/1	0.70	0.44	-	99,99,99,99	0
56	MG	13	1709	1/1	0.94	0.28	-	141,141,141,141	0
56	MG	39	302	1/1	0.78	0.30	-	118,118,118,118	0
56	MG	1H	3285	1/1	0.84	0.39	-	88,88,88,88	0
56	MG	13	1649	1/1	0.97	0.27	-	85,85,85,85	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
56	MG	14	3416	1/1	0.71	0.43	-	126,126,126,126	0
56	MG	13	1710	1/1	0.45	0.30	-	98,98,98,98	0
56	MG	14	3160	1/1	0.92	0.32	-	85,85,85,85	0
56	MG	14	3261	1/1	0.82	0.34	-	83,83,83,83	0
56	MG	1H	3410	1/1	0.93	0.12	-	73,73,73,73	0
56	MG	1H	3041	1/1	0.92	0.33	-	67,67,67,67	0
56	MG	14	3141	1/1	0.94	0.19	-	80,80,80,80	0
56	MG	1H	3071	1/1	0.90	0.12	-	94,94,94,94	0
56	MG	1H	3141	1/1	0.85	0.44	-	89,89,89,89	0
56	MG	1H	3102	1/1	0.86	0.28	-	67,67,67,67	0
56	MG	1H	3297	1/1	0.92	0.10	-	86,86,86,86	0
56	MG	1H	3437	1/1	0.93	0.06	-	109,109,109,109	0
56	MG	1H	3441	1/1	0.94	0.10	-	122,122,122,122	0
56	MG	14	3385	1/1	0.99	0.12	-	73,73,73,73	0
56	MG	14	3364	1/1	0.95	0.07	-	87,87,87,87	0
56	MG	14	3413	1/1	0.87	0.31	-	110,110,110,110	0
56	MG	14	3182	1/1	0.70	0.27	-	98,98,98,98	0
56	MG	13	1678	1/1	0.94	0.09	-	111,111,111,111	0
56	MG	1G	1651	1/1	0.97	0.20	-	138,138,138,138	0
56	MG	1H	3002	1/1	0.93	0.23	-	76,76,76,76	0
56	MG	14	3419	1/1	0.90	0.25	-	112,112,112,112	0
56	MG	13	1623	1/1	0.96	0.15	-	84,84,84,84	0
56	MG	1H	3408	1/1	0.95	0.09	-	87,87,87,87	0
56	MG	14	3370	1/1	0.82	0.12	-	110,110,110,110	0
56	MG	1H	3248	1/1	0.93	0.23	-	76,76,76,76	0
56	MG	13	1742	1/1	0.98	0.10	-	82,82,82,82	0
56	MG	13	1704	1/1	0.85	0.69	-	118,118,118,118	0
56	MG	1H	3045	1/1	0.98	0.20	-	63,63,63,63	0
56	MG	1J	205	1/1	0.86	0.23	-	101,101,101,101	0
56	MG	1H	3099	1/1	0.95	0.33	-	69,69,69,69	0
56	MG	2K	106	1/1	0.82	0.42	-	115,115,115,115	0
56	MG	14	3415	1/1	0.90	0.40	-	107,107,107,107	0
56	MG	1H	3131	1/1	0.97	0.17	-	56,56,56,56	0
56	MG	1H	3472	1/1	0.99	0.08	-	98,98,98,98	0
56	MG	1H	3027	1/1	0.78	0.32	-	96,96,96,96	0
56	MG	14	3011	1/1	0.96	0.20	-	61,61,61,61	0
56	MG	14	3265	1/1	0.73	0.29	-	88,88,88,88	0
56	MG	1H	3266	1/1	0.72	0.58	-	95,95,95,95	0
56	MG	1H	3475	1/1	0.94	0.18	-	76,76,76,76	0
56	MG	14	3080	1/1	0.79	0.41	-	88,88,88,88	0
56	MG	13	1726	1/1	0.95	0.12	-	90,90,90,90	0
56	MG	1H	3228	1/1	0.93	0.55	-	94,94,94,94	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
56	MG	16	204	1/1	0.91	0.40	-	99,99,99,99	0
56	MG	1H	3289	1/1	0.96	0.22	-	93,93,93,93	0
56	MG	1H	3078	1/1	0.97	0.23	-	70,70,70,70	0
56	MG	14	3318	1/1	0.76	0.19	-	99,99,99,99	0
56	MG	1H	3432	1/1	0.92	0.06	-	113,113,113,113	0
56	MG	1H	3328	1/1	0.64	0.55	-	87,87,87,87	0
56	MG	14	3197	1/1	0.86	0.50	-	119,119,119,119	0
56	MG	1H	3152	1/1	0.76	0.30	-	93,93,93,93	0
56	MG	14	3406	1/1	0.94	0.07	-	99,99,99,99	0
56	MG	1H	3355	1/1	0.98	0.18	-	64,64,64,64	0
56	MG	14	3184	1/1	0.93	0.34	-	94,94,94,94	0
56	MG	1H	3464	1/1	0.68	0.14	-	133,133,133,133	0
56	MG	1G	1671	1/1	0.87	0.40	-	120,120,120,120	0
56	MG	13	1635	1/1	0.95	0.14	-	88,88,88,88	0
56	MG	1H	3275	1/1	0.95	0.30	-	97,97,97,97	0
56	MG	16	208	1/1	0.95	0.55	-	94,94,94,94	0
56	MG	14	3162	1/1	0.82	0.19	-	88,88,88,88	0
56	MG	1H	3479	1/1	0.81	0.16	-	126,126,126,126	0
56	MG	14	3304	1/1	0.91	0.33	-	121,121,121,121	0
56	MG	1H	3257	1/1	0.94	0.33	-	80,80,80,80	0
56	MG	13	1637	1/1	0.91	0.25	-	96,96,96,96	0
56	MG	14	3090	1/1	0.96	0.20	-	89,89,89,89	0
56	MG	1H	3127	1/1	0.81	0.29	-	92,92,92,92	0
56	MG	1H	3448	1/1	0.91	0.15	-	97,97,97,97	0
56	MG	1H	3189	1/1	0.77	0.30	-	81,81,81,81	0
56	MG	14	3242	1/1	0.96	0.16	-	94,94,94,94	0
56	MG	14	3279	1/1	0.84	0.25	-	87,87,87,87	0
56	MG	1H	3401	1/1	0.97	0.16	-	87,87,87,87	0
56	MG	14	3038	1/1	0.99	0.16	-	72,72,72,72	0
56	MG	1H	3176	1/1	0.95	0.30	-	91,91,91,91	0
56	MG	1H	3272	1/1	0.65	0.27	-	96,96,96,96	0
56	MG	14	3247	1/1	0.85	0.14	-	75,75,75,75	0
56	MG	1H	3269	1/1	0.96	0.52	-	90,90,90,90	0
56	MG	1H	3117	1/1	0.96	0.29	-	87,87,87,87	0
56	MG	1H	3327	1/1	0.94	0.39	-	91,91,91,91	0
56	MG	14	3259	1/1	0.91	0.15	-	100,100,100,100	0
56	MG	13	1711	1/1	0.92	0.63	-	126,126,126,126	0
56	MG	16	211	1/1	0.94	0.40	-	79,79,79,79	0
56	MG	1H	3029	1/1	0.63	0.28	-	110,110,110,110	0
56	MG	1H	3457	1/1	0.89	0.05	-	113,113,113,113	0
56	MG	1H	3461	1/1	0.93	0.08	-	110,110,110,110	0
56	MG	14	3097	1/1	0.97	0.12	-	69,69,69,69	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
56	MG	1H	3270	1/1	0.62	0.48	-	98,98,98,98	0
56	MG	14	3296	1/1	0.89	0.20	-	107,107,107,107	0
56	MG	1H	3115	1/1	0.97	0.34	-	95,95,95,95	0
56	MG	13	1725	1/1	0.97	0.11	-	102,102,102,102	0
56	MG	14	3426	1/1	0.97	0.24	-	67,67,67,67	0
56	MG	1H	3326	1/1	0.69	0.54	-	117,117,117,117	0
56	MG	1H	3068	1/1	0.94	0.27	-	98,98,98,98	0
56	MG	14	3081	1/1	0.90	0.31	-	90,90,90,90	0
56	MG	14	3186	1/1	0.93	0.21	-	86,86,86,86	0
56	MG	13	1639	1/1	0.70	0.19	-	109,109,109,109	0
56	MG	1G	1633	1/1	0.49	0.32	-	93,93,93,93	0
56	MG	13	1733	1/1	0.98	0.10	-	102,102,102,102	0
56	MG	14	3122	1/1	0.98	0.19	-	68,68,68,68	0
56	MG	16	212	1/1	0.96	0.43	-	99,99,99,99	0
56	MG	14	3299	1/1	0.93	0.58	-	84,84,84,84	0
56	MG	1G	1630	1/1	0.95	0.47	-	116,116,116,116	0
56	MG	14	3398	1/1	0.87	0.06	-	127,127,127,127	0
56	MG	1H	3292	1/1	0.81	0.23	-	72,72,72,72	0
56	MG	1H	3232	1/1	0.82	0.32	-	80,80,80,80	0
56	MG	13	1718	1/1	0.73	0.41	-	96,96,96,96	0
56	MG	14	3057	1/1	0.96	0.15	-	91,91,91,91	0
56	MG	13	1735	1/1	0.97	0.07	-	96,96,96,96	0
56	MG	1G	1624	1/1	0.97	0.17	-	93,93,93,93	0
56	MG	1H	3094	1/1	0.94	0.39	-	79,79,79,79	0
56	MG	14	3381	1/1	0.87	0.22	-	101,101,101,101	0
56	MG	13	1701	1/1	0.93	0.57	-	103,103,103,103	0
56	MG	1H	3387	1/1	0.96	0.17	-	70,70,70,70	0
56	MG	1H	3165	1/1	0.81	0.42	-	95,95,95,95	0
56	MG	1H	3075	1/1	0.96	0.08	-	59,59,59,59	0
56	MG	14	3032	1/1	0.98	0.20	-	52,52,52,52	0
56	MG	14	3102	1/1	0.92	0.36	-	93,93,93,93	0
56	MG	14	3176	1/1	0.79	0.39	-	101,101,101,101	0
56	MG	13	1628	1/1	0.96	0.35	-	102,102,102,102	0
56	MG	1H	3017	1/1	0.99	0.23	-	79,79,79,79	0
56	MG	14	3392	1/1	0.99	0.08	-	70,70,70,70	0
56	MG	14	3228	1/1	0.96	0.20	-	86,86,86,86	0
56	MG	14	3065	1/1	0.82	0.20	-	87,87,87,87	0
56	MG	14	3001	1/1	0.95	0.12	-	66,66,66,66	0
56	MG	1G	1639	1/1	0.86	0.17	-	89,89,89,89	0
56	MG	1H	3459	1/1	0.97	0.13	-	81,81,81,81	0
56	MG	1G	1626	1/1	0.97	0.27	-	102,102,102,102	0
56	MG	1H	3359	1/1	0.93	0.11	-	83,83,83,83	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
56	MG	1G	1634	1/1	0.87	0.48	-	120,120,120,120	0
56	MG	14	3157	1/1	0.94	0.31	-	76,76,76,76	0
56	MG	14	3071	1/1	0.94	0.12	-	94,94,94,94	0
56	MG	14	3328	1/1	0.87	0.12	-	111,111,111,111	0
56	MG	1H	3300	1/1	0.81	0.35	-	87,87,87,87	0
56	MG	1G	1657	1/1	0.79	0.25	-	106,106,106,106	0
56	MG	1H	3249	1/1	0.80	0.41	-	97,97,97,97	0
56	MG	14	3329	1/1	0.58	0.21	-	120,120,120,120	0
56	MG	13	1707	1/1	0.90	0.25	-	104,104,104,104	0
56	MG	14	3237	1/1	0.77	0.24	-	95,95,95,95	0
56	MG	1G	1641	1/1	0.88	0.29	-	90,90,90,90	0
56	MG	1H	3500	1/1	0.77	0.20	-	122,122,122,122	0
56	MG	14	3289	1/1	0.89	0.18	-	86,86,86,86	0
56	MG	14	3371	1/1	0.89	0.21	-	101,101,101,101	0
56	MG	1H	3177	1/1	0.84	0.44	-	98,98,98,98	0
56	MG	16	215	1/1	0.81	0.08	-	113,113,113,113	0
56	MG	14	3418	1/1	0.74	0.07	-	130,130,130,130	0
56	MG	1G	1681	1/1	0.94	0.18	-	133,133,133,133	0
56	MG	13	1619	1/1	0.93	0.58	-	103,103,103,103	0
56	MG	1G	1685	1/1	0.98	0.08	-	112,112,112,112	0
56	MG	1G	1603	1/1	0.86	0.17	-	118,118,118,118	0
56	MG	4K	101	1/1	0.94	0.19	-	90,90,90,90	0
56	MG	1H	3171	1/1	0.97	0.51	-	75,75,75,75	0
56	MG	14	3420	1/1	0.48	0.07	-	130,130,130,130	0
56	MG	1H	3261	1/1	0.95	0.48	-	99,99,99,99	0
56	MG	1H	3051	1/1	0.93	0.24	-	53,53,53,53	0
56	MG	1H	3490	1/1	0.83	0.13	-	125,125,125,125	0
56	MG	14	3290	1/1	0.94	0.26	-	113,113,113,113	0
56	MG	1H	3243	1/1	0.60	0.36	-	93,93,93,93	0
56	MG	1G	1621	1/1	0.98	0.11	-	88,88,88,88	0
56	MG	1H	3147	1/1	0.77	0.31	-	94,94,94,94	0
56	MG	1H	3342	1/1	0.84	0.17	-	92,92,92,92	0
56	MG	13	1741	1/1	0.94	0.08	-	118,118,118,118	0
56	MG	1G	1656	1/1	0.90	0.30	-	103,103,103,103	0
56	MG	14	3323	1/1	0.94	0.10	-	101,101,101,101	0
56	MG	14	3078	1/1	0.93	0.11	-	97,97,97,97	0
56	MG	13	1697	1/1	0.47	0.26	-	97,97,97,97	0
56	MG	14	3362	1/1	0.94	0.17	-	101,101,101,101	0
56	MG	1H	3168	1/1	0.83	0.30	-	83,83,83,83	0
56	MG	1H	3294	1/1	0.83	0.46	-	100,100,100,100	0
56	MG	13	1640	1/1	0.70	0.23	-	94,94,94,94	0
56	MG	1H	3256	1/1	0.85	0.26	-	88,88,88,88	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
56	MG	1G	1673	1/1	0.81	0.14	-	108,108,108,108	0
56	MG	14	3002	1/1	0.97	0.24	-	70,70,70,70	0
56	MG	1H	3425	1/1	0.94	0.06	-	95,95,95,95	0
56	MG	1H	3154	1/1	0.54	0.22	-	91,91,91,91	0
56	MG	14	3021	1/1	0.95	0.40	-	79,79,79,79	0
56	MG	13	1614	1/1	0.97	0.16	-	98,98,98,98	0
56	MG	1H	3463	1/1	0.95	0.04	-	113,113,113,113	0
56	MG	14	3042	1/1	0.94	0.21	-	87,87,87,87	0
56	MG	14	3156	1/1	0.93	0.18	-	99,99,99,99	0
56	MG	14	3142	1/1	0.84	0.46	-	100,100,100,100	0
56	MG	13	1703	1/1	0.92	0.37	-	96,96,96,96	0
56	MG	1H	3444	1/1	0.95	0.29	-	102,102,102,102	0
56	MG	1H	3331	1/1	0.70	0.83	-	106,106,106,106	0
56	MG	35	201	1/1	0.89	0.21	-	91,91,91,91	0
56	MG	1H	3430	1/1	0.90	0.10	-	116,116,116,116	0
56	MG	14	3271	1/1	0.91	0.12	-	87,87,87,87	0
56	MG	1H	3204	1/1	0.88	0.75	-	104,104,104,104	0
56	MG	1H	3213	1/1	0.95	0.27	-	113,113,113,113	0
56	MG	1H	3162	1/1	0.92	0.24	-	74,74,74,74	0
56	MG	14	3045	1/1	0.96	0.17	-	70,70,70,70	0
56	MG	1H	3181	1/1	0.91	0.34	-	74,74,74,74	0
56	MG	1H	3046	1/1	0.96	0.30	-	72,72,72,72	0
56	MG	14	3409	1/1	0.92	0.07	-	114,114,114,114	0
56	MG	13	1731	1/1	0.94	0.06	-	91,91,91,91	0
56	MG	1H	3150	1/1	0.93	0.36	-	86,86,86,86	0
56	MG	1H	3104	1/1	0.89	0.22	-	78,78,78,78	0
56	MG	13	1737	1/1	0.98	0.14	-	72,72,72,72	0
56	MG	14	3340	1/1	0.95	0.34	-	91,91,91,91	0
56	MG	1H	3287	1/1	0.81	0.42	-	92,92,92,92	0
56	MG	1H	3340	1/1	0.73	0.36	-	93,93,93,93	0
56	MG	1H	3107	1/1	0.81	0.28	-	85,85,85,85	0
56	MG	14	3116	1/1	0.97	0.21	-	73,73,73,73	0
56	MG	1G	1640	1/1	0.79	0.49	-	106,106,106,106	0
56	MG	1H	3382	1/1	0.99	0.10	-	58,58,58,58	0
56	MG	1H	3133	1/1	0.91	0.14	-	84,84,84,84	0
56	MG	14	3074	1/1	0.95	0.13	-	79,79,79,79	0
56	MG	1H	3314	1/1	0.81	0.37	-	76,76,76,76	0
56	MG	14	3394	1/1	0.92	0.07	-	93,93,93,93	0
56	MG	14	3314	1/1	0.94	0.34	-	102,102,102,102	0
56	MG	1H	3014	1/1	0.98	0.42	-	57,57,57,57	0
56	MG	14	3284	1/1	0.57	0.30	-	100,100,100,100	0
56	MG	13	1659	1/1	0.97	0.31	-	109,109,109,109	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
56	MG	1H	3066	1/1	0.95	0.27	-	71,71,71,71	0
56	MG	14	3223	1/1	0.97	0.82	-	81,81,81,81	0
56	MG	1G	1620	1/1	0.73	0.20	-	107,107,107,107	0
56	MG	1H	3164	1/1	0.95	0.43	-	102,102,102,102	0
56	MG	13	1613	1/1	0.94	0.18	-	96,96,96,96	0
56	MG	13	1685	1/1	0.77	0.17	-	112,112,112,112	0
56	MG	1H	3434	1/1	0.86	0.19	-	130,130,130,130	0
56	MG	14	3015	1/1	0.97	0.08	-	81,81,81,81	0
56	MG	1H	3449	1/1	0.97	0.15	-	78,78,78,78	0
56	MG	14	3232	1/1	0.74	0.29	-	103,103,103,103	0
56	MG	1H	3307	1/1	0.78	0.28	-	99,99,99,99	0
56	MG	13	1684	1/1	0.82	0.33	-	106,106,106,106	0
56	MG	1H	3477	1/1	0.83	0.10	-	108,108,108,108	0
56	MG	1H	3378	1/1	0.93	0.09	-	94,94,94,94	0
56	MG	14	3231	1/1	0.89	0.21	-	93,93,93,93	0
56	MG	14	3139	1/1	0.95	0.19	-	67,67,67,67	0
56	MG	1H	3097	1/1	0.92	0.20	-	89,89,89,89	0
56	MG	1H	3197	1/1	0.96	0.41	-	89,89,89,89	0
56	MG	14	3269	1/1	0.90	0.31	-	102,102,102,102	0
56	MG	14	3307	1/1	0.52	0.25	-	87,87,87,87	0
56	MG	1H	3486	1/1	0.96	0.07	-	98,98,98,98	0
56	MG	14	3293	1/1	0.81	0.27	-	95,95,95,95	0
56	MG	14	3229	1/1	0.93	0.19	-	83,83,83,83	0
56	MG	1H	3123	1/1	0.99	0.24	-	68,68,68,68	0
56	MG	2L	102	1/1	0.70	0.26	-	89,89,89,89	0
56	MG	1H	3169	1/1	0.75	0.25	-	85,85,85,85	0
56	MG	1H	3114	1/1	0.96	0.42	-	82,82,82,82	0
56	MG	1H	3447	1/1	0.85	0.17	-	105,105,105,105	0
56	MG	14	3155	1/1	0.97	0.16	-	76,76,76,76	0
56	MG	14	3395	1/1	0.91	0.09	-	96,96,96,96	0
56	MG	14	3048	1/1	0.98	0.20	-	73,73,73,73	0
56	MG	1H	3445	1/1	0.97	0.12	-	83,83,83,83	0
56	MG	14	3368	1/1	0.76	0.12	-	117,117,117,117	0
56	MG	1H	3048	1/1	0.98	0.21	-	65,65,65,65	0
56	MG	14	3064	1/1	0.77	0.26	-	97,97,97,97	0
56	MG	1H	3320	1/1	0.92	0.73	-	106,106,106,106	0
56	MG	1H	3156	1/1	0.96	0.36	-	110,110,110,110	0
56	MG	1H	3128	1/1	0.83	0.21	-	88,88,88,88	0
56	MG	1H	3193	1/1	0.83	0.36	-	95,95,95,95	0
56	MG	1H	3242	1/1	0.81	0.52	-	105,105,105,105	0
56	MG	1H	3091	1/1	0.97	0.36	-	70,70,70,70	0
56	MG	1H	3043	1/1	0.98	0.27	-	72,72,72,72	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
56	MG	1H	3190	1/1	0.84	0.17	-	77,77,77,77	0
56	MG	1H	3504	1/1	0.75	0.11	-	138,138,138,138	0
56	MG	2K	107	1/1	0.91	0.41	-	106,106,106,106	0
56	MG	1H	3151	1/1	0.96	0.17	-	110,110,110,110	0
56	MG	1H	3033	1/1	0.96	0.59	-	100,100,100,100	0
56	MG	1H	3108	1/1	0.90	0.22	-	86,86,86,86	0
56	MG	14	3327	1/1	0.98	0.43	-	117,117,117,117	0
56	MG	14	3270	1/1	0.80	0.20	-	107,107,107,107	0
56	MG	14	3243	1/1	0.92	0.10	-	91,91,91,91	0
56	MG	13	1651	1/1	0.71	0.23	-	100,100,100,100	0
56	MG	13	1654	1/1	0.88	0.28	-	123,123,123,123	0
56	MG	1G	1687	1/1	0.84	0.30	-	135,135,135,135	0
56	MG	13	1609	1/1	0.96	0.19	-	83,83,83,83	0
56	MG	14	3180	1/1	0.95	0.31	-	83,83,83,83	0
56	MG	1G	1636	1/1	0.88	0.08	-	115,115,115,115	0
56	MG	14	3163	1/1	0.65	0.18	-	105,105,105,105	0
56	MG	13	1674	1/1	0.89	0.23	-	118,118,118,118	0
56	MG	2L	103	1/1	0.84	0.28	-	115,115,115,115	0
56	MG	1H	3301	1/1	0.94	0.24	-	79,79,79,79	0
56	MG	14	3363	1/1	0.88	0.06	-	117,117,117,117	0
56	MG	13	1708	1/1	0.61	0.34	-	119,119,119,119	0
56	MG	1H	3250	1/1	0.59	0.21	-	104,104,104,104	0
56	MG	1H	3220	1/1	0.94	0.29	-	124,124,124,124	0
56	MG	1H	3098	1/1	0.92	0.23	-	72,72,72,72	0
56	MG	14	3172	1/1	0.96	0.21	-	107,107,107,107	0
56	MG	1H	3259	1/1	0.98	0.52	-	105,105,105,105	0
56	MG	1H	3121	1/1	0.94	0.21	-	80,80,80,80	0

6.5 Other polymers [i](#)

There are no such residues in this entry.