



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 10:18 PM GMT

PDB ID : 4V83  
Title : Crystal structure of a complex containing domain 3 from the PSIV IGR IRES RNA bound to the 70S ribosome.  
Authors : Zhu, J.; Korostelev, A.; Costantino, D.; Noller, H.F.; Kieft, J.S.  
Deposited on : 2010-12-13  
Resolution : 3.50 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](mailto: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.

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

MolProbity : 4.02b-467  
Mogul : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
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) : trunk26865

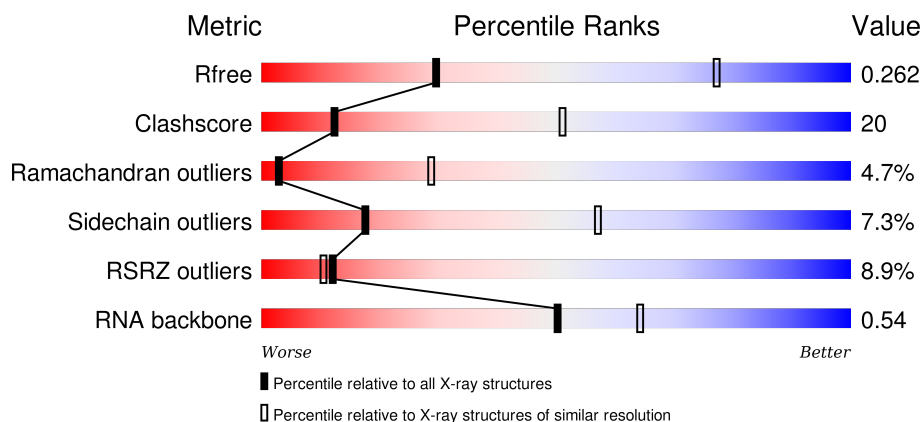
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 3.50 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	1051 (3.60-3.40)
Clashscore	102246	1157 (3.60-3.40)
Ramachandran outliers	100387	1120 (3.60-3.40)
Sidechain outliers	100360	1121 (3.60-3.40)
RSRZ outliers	91569	1058 (3.60-3.40)
RNA backbone	2183	1050 (4.20-2.80)

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  $\geq 3$ , 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	AA	1506	<div> <div>4%</div> <div>47%</div> <div>44%</div> <div>7%</div> <div>.</div> </div>
1	CA	1506	<div> <div>7%</div> <div>47%</div> <div>44%</div> <div>7%</div> <div>.</div> </div>
2	AB	234	<div> <div>34%</div> <div>50%</div> <div>45%</div> <div>5%</div> </div>
2	CB	234	<div> <div>31%</div> <div>51%</div> <div>44%</div> <div>.</div> </div>

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Mol	Chain	Length	Quality of chain
3	AC	206	
3	CC	206	
4	AD	208	
4	CD	208	
5	AE	151	
5	CE	151	
6	AF	101	
6	CF	101	
7	AG	155	
7	CG	155	
8	AH	138	
8	CH	138	
9	AI	127	
9	CI	127	
10	AJ	98	
10	CJ	98	
11	AK	119	
11	CK	119	
12	AL	124	
12	CL	124	
13	AM	116	
13	CM	116	
14	AN	60	
14	CN	60	
15	AO	88	

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Mol	Chain	Length	Quality of chain
15	CO	88	
16	AP	83	
16	CP	83	
17	AQ	99	
17	CQ	99	
18	AR	70	
18	CR	70	
19	AS	78	
19	CS	78	
20	AT	99	
20	CT	99	
21	AU	24	
21	CU	24	
22	AV	35	
22	CV	35	
23	BA	2879	
23	DA	2879	
24	BB	119	
24	DB	119	
25	BC	271	
25	DC	271	
26	BD	204	
26	DD	204	
27	BE	202	
27	DE	202	

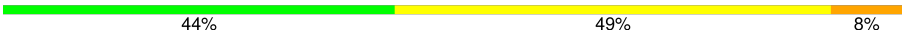
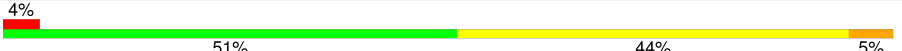
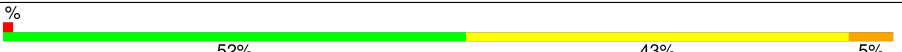
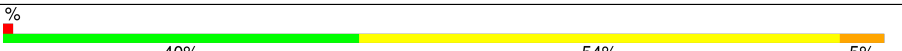
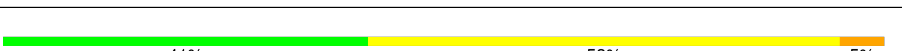
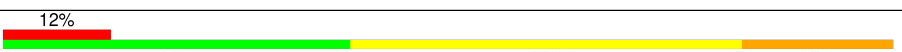
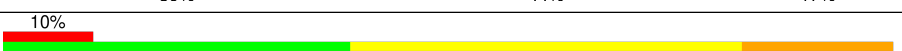

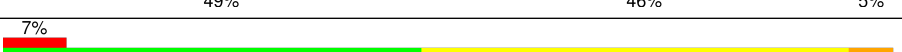

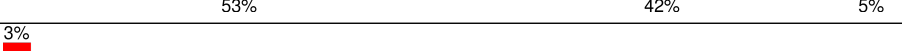
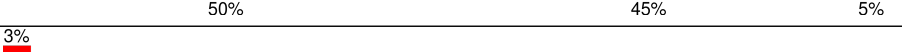


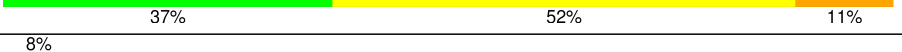
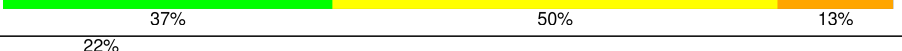



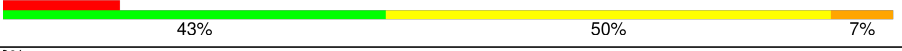
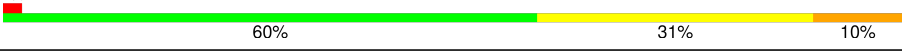


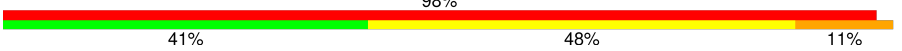

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Mol	Chain	Length	Quality of chain
28	BF	181	
28	DF	181	
29	BG	159	
29	DG	159	
30	BH	145	
30	DH	145	
31	BI	65	
31	DI	65	
32	BJ	137	
32	DJ	137	
33	BK	122	
33	DK	122	
34	BL	146	
34	DL	146	
35	BM	136	
35	DM	136	
36	BN	117	
36	DN	117	
37	BO	98	
37	DO	98	
38	BP	137	
38	DP	137	
39	BQ	116	
39	DQ	116	
40	BR	101	

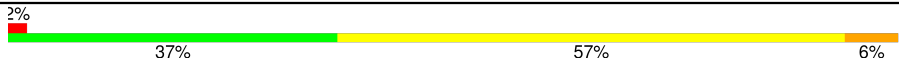

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Mol	Chain	Length	Quality of chain
40	DR	101	
41	BS	112	
41	DS	112	
42	BT	92	
42	DT	92	
43	BU	100	
43	DU	100	
44	BV	188	
44	DV	188	
45	BW	76	
45	DW	76	
46	BX	88	
46	DX	88	
47	BY	62	
47	DY	62	
48	BZ	59	
48	DZ	59	
49	B1	30	
49	D1	30	
50	B2	52	
50	D2	52	
51	B3	44	
51	D3	44	
52	B4	48	
52	D4	48	

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Mol	Chain	Length	Quality of chain
53	B5	63	
53	D5	63	

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
54	MG	AA	1603	-	-	-	X
54	MG	AA	1606	-	-	-	X
54	MG	AA	1609	-	-	-	X
54	MG	AA	1610	-	-	-	X
54	MG	AA	1613	-	-	-	X
54	MG	AA	1615	-	-	-	X
54	MG	AA	1620	-	-	-	X
54	MG	AA	1625	-	-	-	X
54	MG	AA	1626	-	-	-	X
54	MG	AA	1628	-	-	-	X
54	MG	AA	1630	-	-	-	X
54	MG	AA	1631	-	-	-	X
54	MG	AA	1632	-	-	-	X
54	MG	AA	1635	-	-	-	X
54	MG	AA	1641	-	-	-	X
54	MG	AA	1647	-	-	-	X
54	MG	AA	1651	-	-	-	X
54	MG	AA	1664	-	-	-	X
54	MG	AA	1665	-	-	-	X
54	MG	AA	1667	-	-	-	X
54	MG	AA	1681	-	-	-	X
54	MG	AA	1685	-	-	-	X
54	MG	AA	1711	-	-	-	X
54	MG	AA	1720	-	-	-	X
54	MG	AA	1724	-	-	-	X
54	MG	AA	1729	-	-	-	X
54	MG	AA	1760	-	-	-	X
54	MG	AA	1778	-	-	-	X
54	MG	AA	1784	-	-	-	X
54	MG	AA	1800	-	-	-	X
54	MG	AA	1820	-	-	-	X
54	MG	AA	1828	-	-	-	X
54	MG	AA	1846	-	-	-	X

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
54	MG	AA	1860	-	-	-	X
54	MG	AA	1870	-	-	-	X
54	MG	AD	303	-	-	-	X
54	MG	AT	202	-	-	-	X
54	MG	B5	101	-	-	-	X
54	MG	BA	2902	-	-	-	X
54	MG	BA	2903	-	-	-	X
54	MG	BA	2906	-	-	-	X
54	MG	BA	2907	-	-	-	X
54	MG	BA	2909	-	-	-	X
54	MG	BA	2911	-	-	-	X
54	MG	BA	2912	-	-	-	X
54	MG	BA	2913	-	-	-	X
54	MG	BA	2914	-	-	-	X
54	MG	BA	2916	-	-	-	X
54	MG	BA	2919	-	-	-	X
54	MG	BA	2920	-	-	-	X
54	MG	BA	2922	-	-	-	X
54	MG	BA	2923	-	-	-	X
54	MG	BA	2924	-	-	-	X
54	MG	BA	2926	-	-	-	X
54	MG	BA	2929	-	-	-	X
54	MG	BA	2931	-	-	-	X
54	MG	BA	2934	-	-	-	X
54	MG	BA	2939	-	-	-	X
54	MG	BA	2943	-	-	-	X
54	MG	BA	2945	-	-	-	X
54	MG	BA	2954	-	-	-	X
54	MG	BA	2955	-	-	-	X
54	MG	BA	2957	-	-	-	X
54	MG	BA	2960	-	-	-	X
54	MG	BA	2963	-	-	-	X
54	MG	BA	2968	-	-	-	X
54	MG	BA	2975	-	-	-	X
54	MG	BA	2981	-	-	-	X
54	MG	BA	2990	-	-	-	X
54	MG	BA	2991	-	-	-	X
54	MG	BA	3003	-	-	-	X
54	MG	BA	3004	-	-	-	X
54	MG	BA	3008	-	-	-	X
54	MG	BA	3013	-	-	-	X
54	MG	BA	3019	-	-	-	X

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
54	MG	BA	3020	-	-	-	X
54	MG	BA	3021	-	-	-	X
54	MG	BA	3022	-	-	-	X
54	MG	BA	3023	-	-	-	X
54	MG	BA	3038	-	-	-	X
54	MG	BA	3040	-	-	-	X
54	MG	BA	3047	-	-	-	X
54	MG	BA	3048	-	-	-	X
54	MG	BA	3049	-	-	-	X
54	MG	BA	3051	-	-	-	X
54	MG	BA	3058	-	-	-	X
54	MG	BA	3060	-	-	-	X
54	MG	BA	3064	-	-	-	X
54	MG	BA	3068	-	-	-	X
54	MG	BA	3070	-	-	-	X
54	MG	BA	3072	-	-	-	X
54	MG	BA	3073	-	-	-	X
54	MG	BA	3088	-	-	-	X
54	MG	BA	3098	-	-	-	X
54	MG	BA	3110	-	-	-	X
54	MG	BA	3114	-	-	-	X
54	MG	BA	3122	-	-	-	X
54	MG	BA	3126	-	-	-	X
54	MG	BA	3136	-	-	-	X
54	MG	BA	3150	-	-	-	X
54	MG	BA	3159	-	-	-	X
54	MG	BA	3164	-	-	-	X
54	MG	BA	3169	-	-	-	X
54	MG	BA	3172	-	-	-	X
54	MG	BA	3175	-	-	-	X
54	MG	BA	3178	-	-	-	X
54	MG	BA	3187	-	-	-	X
54	MG	BA	3196	-	-	-	X
54	MG	BA	3227	-	-	-	X
54	MG	BA	3238	-	-	-	X
54	MG	BA	3239	-	-	-	X
54	MG	BA	3243	-	-	-	X
54	MG	BA	3289	-	-	-	X
54	MG	BA	3301	-	-	-	X
54	MG	BA	3327	-	-	-	X
54	MG	BA	3334	-	-	-	X
54	MG	BA	3340	-	-	-	X

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
54	MG	BA	3411	-	-	-	X
54	MG	BA	3425	-	-	-	X
54	MG	BA	3432	-	-	-	X
54	MG	BA	3434	-	-	-	X
54	MG	BA	3438	-	-	-	X
54	MG	BA	3442	-	-	-	X
54	MG	BA	3443	-	-	-	X
54	MG	BA	3451	-	-	-	X
54	MG	BA	3455	-	-	-	X
54	MG	BA	3459	-	-	-	X
54	MG	BA	3461	-	-	-	X
54	MG	BA	3467	-	-	-	X
54	MG	BA	3469	-	-	-	X
54	MG	BA	3471	-	-	-	X
54	MG	BA	3556	-	-	-	X
54	MG	BA	3570	-	-	-	X
54	MG	BA	3575	-	-	-	X
54	MG	BA	3597	-	-	-	X
54	MG	BA	3626	-	-	-	X
54	MG	BA	3634	-	-	-	X
54	MG	BA	3643	-	-	-	X
54	MG	BA	3673	-	-	-	X
54	MG	BB	218	-	-	-	X
54	MG	BC	303	-	-	-	X
54	MG	BE	302	-	-	-	X
54	MG	BM	203	-	-	-	X
54	MG	BX	102	-	-	-	X
54	MG	CA	1601	-	-	-	X
54	MG	CA	1602	-	-	-	X
54	MG	CA	1606	-	-	-	X
54	MG	CA	1619	-	-	-	X
54	MG	CA	1651	-	-	-	X
54	MG	CA	1659	-	-	-	X
54	MG	CA	1665	-	-	-	X
54	MG	CA	1676	-	-	-	X
54	MG	CA	1680	-	-	-	X
54	MG	CA	1688	-	-	-	X
54	MG	CA	1733	-	-	-	X
54	MG	CA	1737	-	-	-	X
54	MG	CA	1760	-	-	-	X
54	MG	CA	1788	-	-	-	X
54	MG	CA	1862	-	-	-	X

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
54	MG	CA	1875	-	-	-	X
54	MG	CA	1881	-	-	-	X
54	MG	CA	1891	-	-	-	X
54	MG	CA	1915	-	-	-	X
54	MG	CA	1943	-	-	-	X
54	MG	CD	304	-	-	-	X
54	MG	CM	201	-	-	-	X
54	MG	CN	102	-	-	-	X
54	MG	DA	2901	-	-	-	X
54	MG	DA	2902	-	-	-	X
54	MG	DA	2904	-	-	-	X
54	MG	DA	2906	-	-	-	X
54	MG	DA	2907	-	-	-	X
54	MG	DA	2908	-	-	-	X
54	MG	DA	2913	-	-	-	X
54	MG	DA	2918	-	-	-	X
54	MG	DA	2923	-	-	-	X
54	MG	DA	2925	-	-	-	X
54	MG	DA	2926	-	-	-	X
54	MG	DA	2927	-	-	-	X
54	MG	DA	2928	-	-	-	X
54	MG	DA	2932	-	-	-	X
54	MG	DA	2934	-	-	-	X
54	MG	DA	2938	-	-	-	X
54	MG	DA	2943	-	-	-	X
54	MG	DA	2947	-	-	-	X
54	MG	DA	2953	-	-	-	X
54	MG	DA	2954	-	-	-	X
54	MG	DA	2957	-	-	-	X
54	MG	DA	2971	-	-	-	X
54	MG	DA	2973	-	-	-	X
54	MG	DA	2975	-	-	-	X
54	MG	DA	2980	-	-	-	X
54	MG	DA	2982	-	-	-	X
54	MG	DA	2985	-	-	-	X
54	MG	DA	2986	-	-	-	X
54	MG	DA	2997	-	-	-	X
54	MG	DA	3005	-	-	-	X
54	MG	DA	3010	-	-	-	X
54	MG	DA	3013	-	-	-	X
54	MG	DA	3014	-	-	-	X
54	MG	DA	3015	-	-	-	X

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
54	MG	DA	3018	-	-	-	X
54	MG	DA	3020	-	-	-	X
54	MG	DA	3024	-	-	-	X
54	MG	DA	3028	-	-	-	X
54	MG	DA	3035	-	-	-	X
54	MG	DA	3038	-	-	-	X
54	MG	DA	3045	-	-	-	X
54	MG	DA	3048	-	-	-	X
54	MG	DA	3052	-	-	-	X
54	MG	DA	3056	-	-	-	X
54	MG	DA	3057	-	-	-	X
54	MG	DA	3059	-	-	-	X
54	MG	DA	3063	-	-	-	X
54	MG	DA	3064	-	-	-	X
54	MG	DA	3065	-	-	-	X
54	MG	DA	3072	-	-	-	X
54	MG	DA	3074	-	-	-	X
54	MG	DA	3080	-	-	-	X
54	MG	DA	3082	-	-	-	X
54	MG	DA	3098	-	-	-	X
54	MG	DA	3103	-	-	-	X
54	MG	DA	3106	-	-	-	X
54	MG	DA	3108	-	-	-	X
54	MG	DA	3121	-	-	-	X
54	MG	DA	3126	-	-	-	X
54	MG	DA	3144	-	-	-	X
54	MG	DA	3152	-	-	-	X
54	MG	DA	3153	-	-	-	X
54	MG	DA	3154	-	-	-	X
54	MG	DA	3158	-	-	-	X
54	MG	DA	3160	-	-	-	X
54	MG	DA	3161	-	-	-	X
54	MG	DA	3172	-	-	-	X
54	MG	DA	3173	-	-	-	X
54	MG	DA	3177	-	-	-	X
54	MG	DA	3178	-	-	-	X
54	MG	DA	3180	-	-	-	X
54	MG	DA	3185	-	-	-	X
54	MG	DA	3186	-	-	-	X
54	MG	DA	3193	-	-	-	X
54	MG	DA	3197	-	-	-	X
54	MG	DA	3226	-	-	-	X

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
54	MG	DA	3227	-	-	-	X
54	MG	DA	3258	-	-	-	X
54	MG	DA	3267	-	-	-	X
54	MG	DA	3268	-	-	-	X
54	MG	DA	3289	-	-	-	X
54	MG	DA	3310	-	-	-	X
54	MG	DA	3311	-	-	-	X
54	MG	DA	3320	-	-	-	X
54	MG	DA	3321	-	-	-	X
54	MG	DA	3329	-	-	-	X
54	MG	DA	3333	-	-	-	X
54	MG	DA	3343	-	-	-	X
54	MG	DA	3353	-	-	-	X
54	MG	DA	3357	-	-	-	X
54	MG	DA	3360	-	-	-	X
54	MG	DA	3412	-	-	-	X
54	MG	DA	3414	-	-	-	X
54	MG	DA	3415	-	-	-	X
54	MG	DA	3431	-	-	-	X
54	MG	DA	3455	-	-	-	X
54	MG	DA	3472	-	-	-	X
54	MG	DA	3506	-	-	-	X
54	MG	DA	3511	-	-	-	X
54	MG	DA	3529	-	-	-	X
54	MG	DA	3534	-	-	-	X
54	MG	DA	3563	-	-	-	X
54	MG	DA	3569	-	-	-	X
54	MG	DA	3590	-	-	-	X
54	MG	DA	3599	-	-	-	X
54	MG	DA	3603	-	-	-	X
54	MG	DA	3605	-	-	-	X
54	MG	DA	3606	-	-	-	X
54	MG	DA	3607	-	-	-	X
54	MG	DA	3622	-	-	-	X
54	MG	DA	3629	-	-	-	X
54	MG	DA	3633	-	-	-	X
54	MG	DA	3634	-	-	-	X
54	MG	DA	3637	-	-	-	X
54	MG	DA	3638	-	-	-	X
54	MG	DA	3646	-	-	-	X
54	MG	DA	3650	-	-	-	X
54	MG	DA	3655	-	-	-	X

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
54	MG	DA	3658	-	-	-	X
54	MG	DA	3667	-	-	-	X
54	MG	DA	3672	-	-	-	X
54	MG	DA	3684	-	-	-	X
54	MG	DA	3689	-	-	-	X
54	MG	DA	3698	-	-	-	X
54	MG	DA	3733	-	-	-	X
54	MG	DA	3736	-	-	-	X
54	MG	DA	3742	-	-	-	X
54	MG	DA	3750	-	-	-	X
54	MG	DA	3771	-	-	-	X
54	MG	DA	3840	-	-	-	X
54	MG	DA	3851	-	-	-	X
54	MG	DA	3859	-	-	-	X
54	MG	DA	3863	-	-	-	X
54	MG	DB	204	-	-	-	X
54	MG	DB	207	-	-	-	X
54	MG	DB	216	-	-	-	X
54	MG	DC	302	-	-	-	X
54	MG	DD	301	-	-	-	X
54	MG	DD	305	-	-	-	X
54	MG	DJ	204	-	-	-	X
54	MG	DJ	205	-	-	-	X
54	MG	DW	101	-	-	-	X
54	MG	DX	101	-	-	-	X

## 2 Entry composition

There are 55 unique types of molecules in this entry. The entry contains 283641 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 ribosomal RNA 16S.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	AA	1506	Total	C	N	O	P	0	0	0
			32372	14409	5999	10459	1505			
1	CA	1506	Total	C	N	O	P	0	0	0
			32372	14409	5999	10459	1505			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	AB	234	Total	C	N	O	S	0	0	0
			1901	1213	341	342	5			
2	CB	234	Total	C	N	O	S	0	0	0
			1901	1213	341	342	5			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	AC	206	Total	C	N	O	S	0	0	0
			1613	1016	314	282	1			
3	CC	206	Total	C	N	O	S	0	0	0
			1613	1016	314	282	1			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	AD	208	Total	C	N	O	S	0	0	0
			1703	1066	339	291	7			
4	CD	208	Total	C	N	O	S	0	0	0
			1703	1066	339	291	7			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	AE	151	Total	C	N	O	S	0	0	0
			1156	729	218	205	4			
5	CE	151	Total	C	N	O	S	0	0	0
			1156	729	218	205	4			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
6	AF	101	Total	C	N	O	S	0	0	0
			843	531	155	154	3			
6	CF	101	Total	C	N	O	S	0	0	0
			843	531	155	154	3			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
7	AG	155	Total	C	N	O	S	0	0	0
			1257	781	252	218	6			
7	CG	155	Total	C	N	O	S	0	0	0
			1257	781	252	218	6			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
8	AH	138	Total	C	N	O	S	0	0	0
			1116	705	215	193	3			
8	CH	138	Total	C	N	O	S	0	0	0
			1116	705	215	193	3			

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

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
9	AI	127	Total	C	N	O	0	0	0
			1011	639	198	174			
9	CI	127	Total	C	N	O	0	0	0
			1011	639	198	174			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
10	AJ	98	Total	C	N	O	S	0	0	0
			795	499	156	139	1			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
10	CJ	98	Total	C	N	O	S	0	0	0
			795	499	156	139	1			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
11	AK	119	Total	C	N	O	S	0	0	0
			885	549	168	165	3			
11	CK	119	Total	C	N	O	S	0	0	0
			885	549	168	165	3			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
12	AL	124	Total	C	N	O	S	0	0	0
			971	611	195	164	1			
12	CL	124	Total	C	N	O	S	0	0	0
			971	611	195	164	1			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
13	AM	116	Total	C	N	O	S	0	0	0
			929	574	191	162	2			
13	CM	116	Total	C	N	O	S	0	0	0
			929	574	191	162	2			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
14	AN	60	Total	C	N	O	S	0	0	0
			492	312	104	72	4			
14	CN	60	Total	C	N	O	S	0	0	0
			492	312	104	72	4			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
15	AO	88	Total	C	N	O	S	0	0	0
			734	459	147	126	2			
15	CO	88	Total	C	N	O	S	0	0	0
			734	459	147	126	2			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
16	AP	83	Total	C	N	O	S	0	0	0
			701	443	139	118	1			
16	CP	83	Total	C	N	O	S	0	0	0
			701	443	139	118	1			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
17	AQ	99	Total	C	N	O	S	0	0	0
			824	528	152	142	2			
17	CQ	99	Total	C	N	O	S	0	0	0
			824	528	152	142	2			

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

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
18	AR	70	Total	C	N	O	0	0	0
			574	367	112	95			
18	CR	70	Total	C	N	O	0	0	0
			574	367	112	95			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
19	AS	78	Total	C	N	O	S	0	0	0
			630	403	114	111	2			
19	CS	78	Total	C	N	O	S	0	0	0
			630	403	114	111	2			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
20	AT	99	Total	C	N	O	S	0	0	0
			762	469	162	129	2			
20	CT	99	Total	C	N	O	S	0	0	0
			762	469	162	129	2			

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

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
21	AU	24	Total	C	N	O	0	0	0
			209	128	50	31			
21	CU	24	Total	C	N	O	0	0	0
			209	128	50	31			

- Molecule 22 is a RNA chain called domain 3 of PSIC IGR IRES RNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
22	AV	35	Total	C	N	O	P	0	0	0
			736	332	128	243	33			
22	CV	35	Total	C	N	O	P	0	0	0
			736	332	128	243	33			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
23	BA	2760	Total	C	N	O	P	0	0	0
			59442	26456	11114	19113	2759			
23	DA	2760	Total	C	N	O	P	0	0	0
			59442	26456	11114	19113	2759			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
BA	1142	U	C	SEE REMARK 999	GB AE017221.1
BA	2825	U	G	SEE REMARK 999	GB AE017221.1
DA	1142	U	C	SEE REMARK 999	GB AE017221.1
DA	2825	U	G	SEE REMARK 999	GB AE017221.1

- Molecule 24 is a RNA chain called 5S ribosomal RNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
24	BB	119	Total	C	N	O	P	0	0	0
			2551	1136	471	826	118			
24	DB	119	Total	C	N	O	P	0	0	0
			2551	1136	471	826	118			

- Molecule 25 is a protein called 50S ribosomal protein L2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
25	BC	271	Total	C	N	O	S	0	0	0
			2105	1329	416	357	3			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
25	DC	271	Total	C	N	O	S	0	0	0
			2105	1329	416	357	3			

- Molecule 26 is a protein called 50S ribosomal protein L3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
26	BD	204	Total	C	N	O	S	0	0	0
			1564	988	299	271	6			
26	DD	204	Total	C	N	O	S	0	0	0
			1564	988	299	271	6			

- Molecule 27 is a protein called 50S ribosomal protein L4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
27	BE	202	Total	C	N	O	S	0	0	0
			1587	1011	297	276	3			
27	DE	202	Total	C	N	O	S	0	0	0
			1587	1011	297	276	3			

- Molecule 28 is a protein called 50S ribosomal protein L5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
28	BF	181	Total	C	N	O	S	0	0	0
			1475	943	268	260	4			
28	DF	181	Total	C	N	O	S	0	0	0
			1475	943	268	260	4			

- Molecule 29 is a protein called 50S ribosomal protein L6.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
29	BG	159	Total	C	N	O	S	0	0	0
			1223	773	228	221	1			
29	DG	159	Total	C	N	O	S	0	0	0
			1223	773	228	221	1			

- Molecule 30 is a protein called 50S ribosomal protein L9.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
30	BH	145	Total	C	N	O	S	0	0	0
			1133	724	200	208	1			
30	DH	145	Total	C	N	O	S	0	0	0
			1133	724	200	208	1			

- Molecule 31 is a protein called 50S ribosomal protein L10.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
31	BI	32	Total	C	N	O	0	0	0
			254	157	49	48			
31	DI	32	Total	C	N	O	0	0	0
			254	157	49	48			

- Molecule 32 is a protein called 50S ribosomal protein L13.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
32	BJ	137	Total	C	N	O	S	0	0	0
			1097	707	205	182	3			
32	DJ	137	Total	C	N	O	S	0	0	0
			1097	707	205	182	3			

- Molecule 33 is a protein called 50S ribosomal protein L14.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
33	BK	122	Total	C	N	O	S	0	0	0
			932	587	171	170	4			
33	DK	122	Total	C	N	O	S	0	0	0
			932	587	171	170	4			

- Molecule 34 is a protein called 50S ribosomal protein L15.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
34	BL	146	Total	C	N	O	S	0	0	0
			1114	692	227	193	2			
34	DL	146	Total	C	N	O	S	0	0	0
			1114	692	227	193	2			

- Molecule 35 is a protein called 50S ribosomal protein L16.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
35	BM	136	Total	C	N	O	S	0	0	0
			1079	688	204	182	5			
35	DM	136	Total	C	N	O	S	0	0	0
			1079	688	204	182	5			

- Molecule 36 is a protein called 50S ribosomal protein L17.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
36	BN	117	Total	C	N	O	0	0	0
			960	599	202	159			
36	DN	117	Total	C	N	O	0	0	0
			960	599	202	159			

- Molecule 37 is a protein called 50S ribosomal protein L18.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
37	BO	98	Total	C	N	O	0	0	0
			771	486	154	131			
37	DO	98	Total	C	N	O	0	0	0
			771	486	154	131			

- Molecule 38 is a protein called 50S ribosomal protein L19.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
38	BP	137	Total	C	N	O	S	0	0	0
			1144	713	234	196	1			
38	DP	137	Total	C	N	O	S	0	0	0
			1144	713	234	196	1			

- Molecule 39 is a protein called 50S ribosomal protein L20.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
39	BQ	116	Total	C	N	O	S	0	0	0
			953	601	201	150	1			
39	DQ	116	Total	C	N	O	S	0	0	0
			953	601	201	150	1			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
BQ	?	-	PHE	DELETION	UNP Q72L76
DQ	?	-	PHE	DELETION	UNP Q72L76

- Molecule 40 is a protein called 50S ribosomal protein L21.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
40	BR	101	Total	C	N	O	S	0	0	0
			779	501	142	135	1			
40	DR	101	Total	C	N	O	S	0	0	0
			779	501	142	135	1			

- Molecule 41 is a protein called 50S ribosomal protein L22.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
41	BS	112	Total	C	N	O	S	0	0	0
			891	560	175	154	2			
41	DS	112	Total	C	N	O	S	0	0	0
			891	560	175	154	2			

- Molecule 42 is a protein called 50S ribosomal protein L23.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
42	BT	92	Total	C	N	O		0	0	0
			726	471	131	124				
42	DT	92	Total	C	N	O		0	0	0
			726	471	131	124				

- Molecule 43 is a protein called 50S ribosomal protein L24.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
43	BU	100	Total	C	N	O	S	0	0	0
			776	500	148	124	4			
43	DU	100	Total	C	N	O	S	0	0	0
			776	500	148	124	4			

- Molecule 44 is a protein called 50S ribosomal protein L25.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
44	BV	188	Total	C	N	O	S	0	0	0
			1492	950	265	275	2			
44	DV	188	Total	C	N	O	S	0	0	0
			1492	950	265	275	2			

- Molecule 45 is a protein called 50S ribosomal protein L27.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
45	BW	76	Total	C	N	O	S	0	0	0
			605	376	126	102	1			
45	DW	76	Total	C	N	O	S	0	0	0
			605	376	126	102	1			

- Molecule 46 is a protein called 50S ribosomal protein L28.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
46	BX	88	Total	C	N	O	0	0	0
			695	435	141	119			
46	DX	88	Total	C	N	O	0	0	0
			695	435	141	119			

- Molecule 47 is a protein called 50S ribosomal protein L29.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
47	BY	62	Total	C	N	O	S	0	0	0
			521	325	102	92	2			
47	DY	62	Total	C	N	O	S	0	0	0
			521	325	102	92	2			

- Molecule 48 is a protein called 50S ribosomal protein L30.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
48	BZ	59	Total	C	N	O	S	0	0	0
			468	298	90	79	1			
48	DZ	59	Total	C	N	O	S	0	0	0
			468	298	90	79	1			

- Molecule 49 is a protein called 50S ribosomal protein L31.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
49	B1	30	Total	C	N	O	S	0	0	0
			226	142	36	44	4			
49	D1	30	Total	C	N	O	S	0	0	0
			226	142	36	44	4			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
50	B2	52	Total	C	N	O	S	0	0	0
			405	255	79	66	5			
50	D2	52	Total	C	N	O	S	0	0	0
			405	255	79	66	5			

- Molecule 51 is a protein called 50S ribosomal protein L33.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
51	B3	44	Total	C	N	O	S	0	0	0
			381	235	77	65	4			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
51	D3	44	Total	C	N	O	S	0	0	0
			381	235	77	65	4			

- Molecule 52 is a protein called 50S ribosomal protein L34.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
52	B4	48	Total	C	N	O	S	0	0	0
			419	257	104	56	2			
52	D4	48	Total	C	N	O	S	0	0	0
			419	257	104	56	2			

- Molecule 53 is a protein called 50S ribosomal protein L35.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
53	B5	63	Total	C	N	O	S	0	0	0
			508	326	101	79	2			
53	D5	63	Total	C	N	O	S	0	0	0
			508	326	101	79	2			

- Molecule 54 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
54	DC	3	Total	Mg	0	0
			3	3		
54	DX	3	Total	Mg	0	0
			3	3		
54	B4	1	Total	Mg	0	0
			1	1		
54	BA	781	Total	Mg	0	0
			781	781		
54	AK	2	Total	Mg	0	0
			2	2		
54	DQ	2	Total	Mg	0	0
			2	2		
54	AB	3	Total	Mg	0	0
			3	3		
54	DF	2	Total	Mg	0	0
			2	2		
54	CV	7	Total	Mg	0	0
			7	7		
54	DL	3	Total	Mg	0	0
			3	3		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
54	BE	2	Total 2	Mg 2	0	0
54	DU	1	Total 1	Mg 1	0	0
54	DY	2	Total 2	Mg 2	0	0
54	AN	1	Total 1	Mg 1	0	0
54	BP	2	Total 2	Mg 2	0	0
54	CN	1	Total 1	Mg 1	0	0
54	BI	1	Total 1	Mg 1	0	0
54	AS	1	Total 1	Mg 1	0	0
54	CA	372	Total 372	Mg 372	0	0
54	B5	1	Total 1	Mg 1	0	0
54	BB	27	Total 27	Mg 27	0	0
54	BT	2	Total 2	Mg 2	0	0
54	DO	1	Total 1	Mg 1	0	0
54	AE	5	Total 5	Mg 5	0	0
54	BM	3	Total 3	Mg 3	0	0
54	CF	2	Total 2	Mg 2	0	0
54	D3	1	Total 1	Mg 1	0	0
54	BF	1	Total 1	Mg 1	0	0
54	AV	8	Total 8	Mg 8	0	0
54	BX	2	Total 2	Mg 2	0	0
54	DA	964	Total 964	Mg 964	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
54	B2	3	Total 3	Mg 3	0	0
54	AA	279	Total 279	Mg 279	0	0
54	CQ	2	Total 2	Mg 2	0	0
54	BJ	3	Total 3	Mg 3	0	0
54	DV	4	Total 4	Mg 4	0	0
54	BC	3	Total 3	Mg 3	0	0
54	AM	1	Total 1	Mg 1	0	0
54	BU	2	Total 2	Mg 2	0	0
54	CC	8	Total 8	Mg 8	0	0
54	AD	3	Total 3	Mg 3	0	0
54	BN	2	Total 2	Mg 2	0	0
54	CM	2	Total 2	Mg 2	0	0
54	DS	2	Total 2	Mg 2	0	0
54	DM	1	Total 1	Mg 1	0	0
54	AI	1	Total 1	Mg 1	0	0
54	BY	3	Total 3	Mg 3	0	0
54	DE	3	Total 3	Mg 3	0	0
54	B3	1	Total 1	Mg 1	0	0
54	DG	3	Total 3	Mg 3	0	0
54	BR	1	Total 1	Mg 1	0	0
54	DK	3	Total 3	Mg 3	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
54	D4	4	Total 4	Mg 4	0	0
54	BK	1	Total 1	Mg 1	0	0
54	CE	5	Total 5	Mg 5	0	0
54	DW	1	Total 1	Mg 1	0	0
54	D5	3	Total 3	Mg 3	0	0
54	DD	6	Total 6	Mg 6	0	0
54	D2	3	Total 3	Mg 3	0	0
54	AL	2	Total 2	Mg 2	0	0
54	BV	3	Total 3	Mg 3	0	0
54	AG	2	Total 2	Mg 2	0	0
54	BO	1	Total 1	Mg 1	0	0
54	AQ	3	Total 3	Mg 3	0	0
54	D1	1	Total 1	Mg 1	0	0
54	DI	1	Total 1	Mg 1	0	0
54	AH	1	Total 1	Mg 1	0	0
54	DJ	6	Total 6	Mg 6	0	0
54	CO	4	Total 4	Mg 4	0	0
54	AC	2	Total 2	Mg 2	0	0
54	BS	1	Total 1	Mg 1	0	0
54	DB	43	Total 43	Mg 43	0	0
54	CB	3	Total 3	Mg 3	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
54	CD	3	Total 3	Mg 3	0	0
54	DN	1	Total 1	Mg 1	0	0
54	AT	2	Total 2	Mg 2	0	0
54	CL	6	Total 6	Mg 6	0	0
54	DP	2	Total 2	Mg 2	0	0
54	AO	3	Total 3	Mg 3	0	0
54	BW	1	Total 1	Mg 1	0	0
54	CG	3	Total 3	Mg 3	0	0
54	CK	1	Total 1	Mg 1	0	0
54	AF	1	Total 1	Mg 1	0	0
54	BH	1	Total 1	Mg 1	0	0

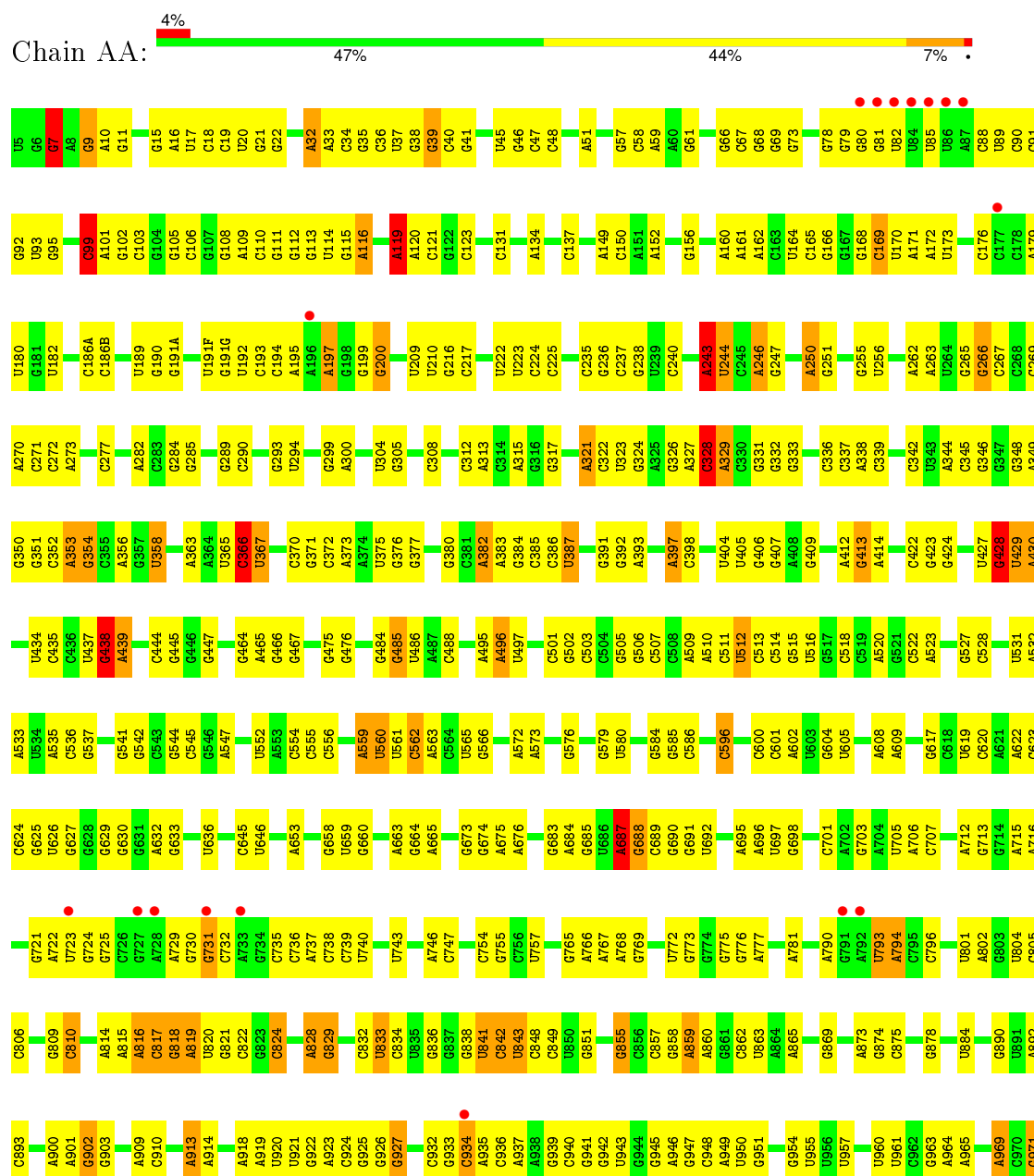
- Molecule 55 is ZINC ION (three-letter code: ZN) (formula: Zn).

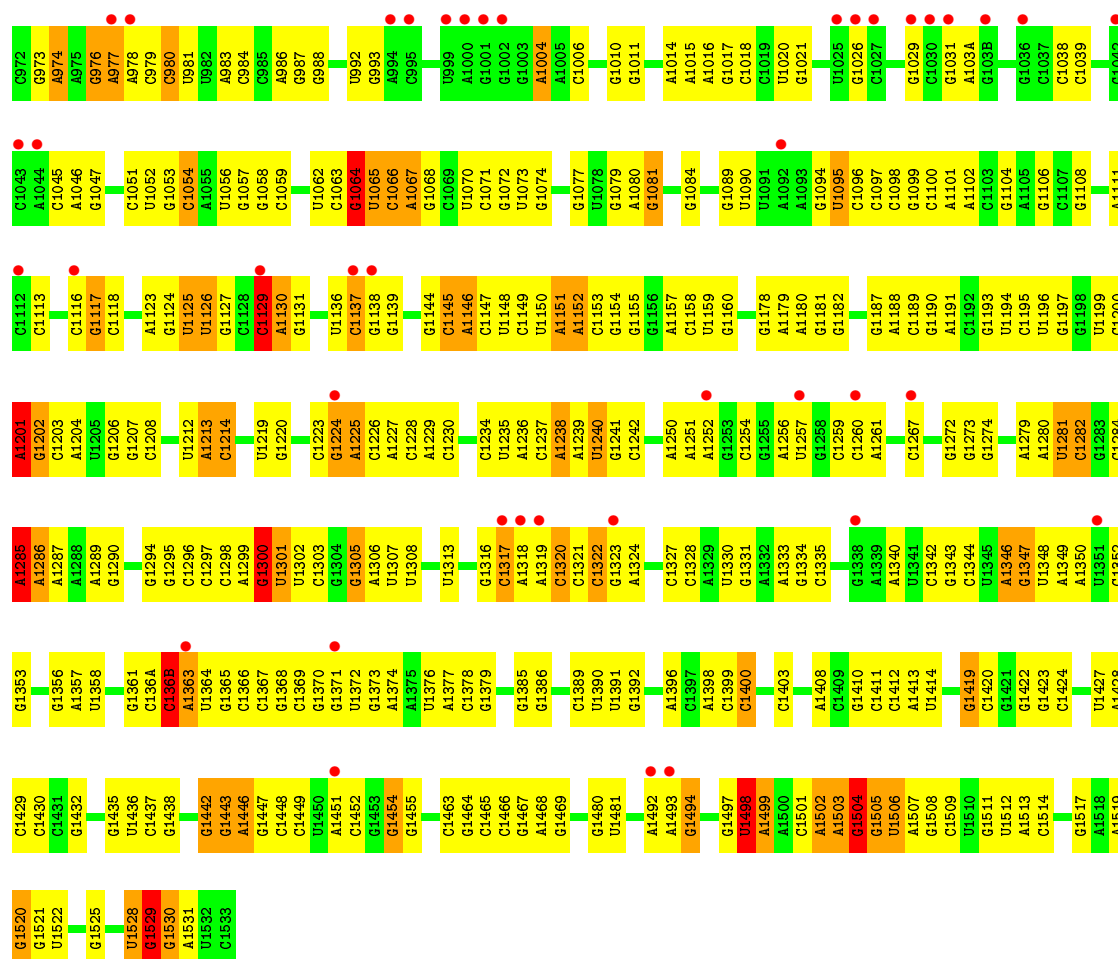
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
55	CN	1	Total 1	Zn 1	0	0
55	AD	1	Total 1	Zn 1	0	0
55	CD	1	Total 1	Zn 1	0	0
55	AN	1	Total 1	Zn 1	0	0

### 3 Residue-property plots

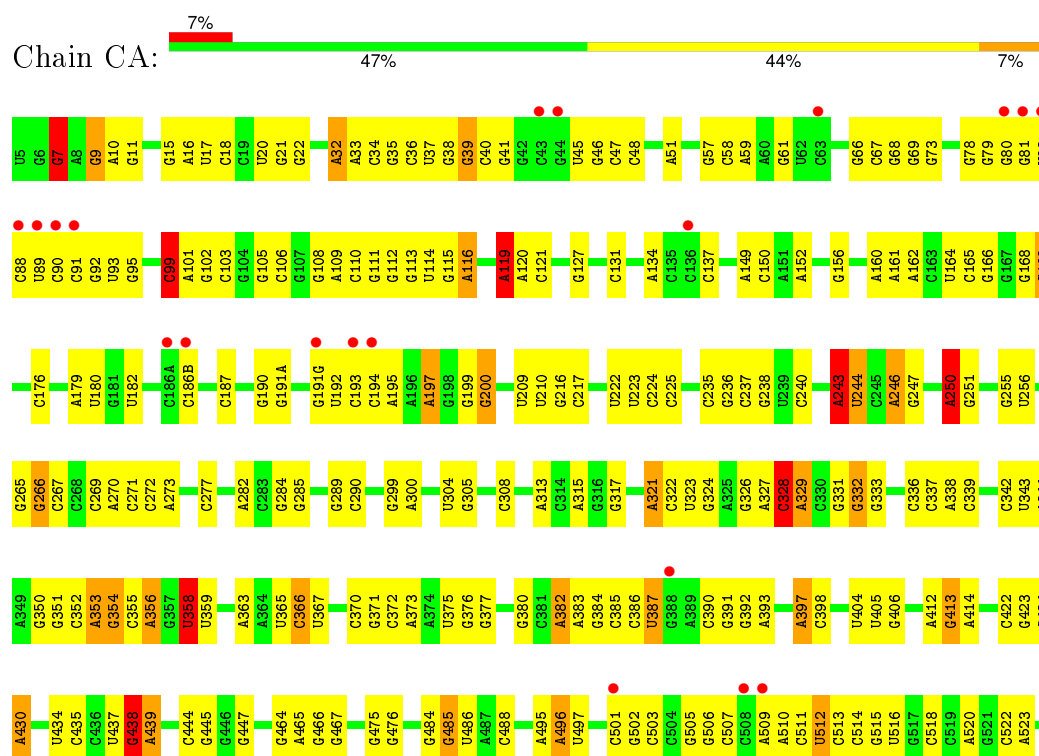
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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 ( $RSRZ > 2$ ). 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.

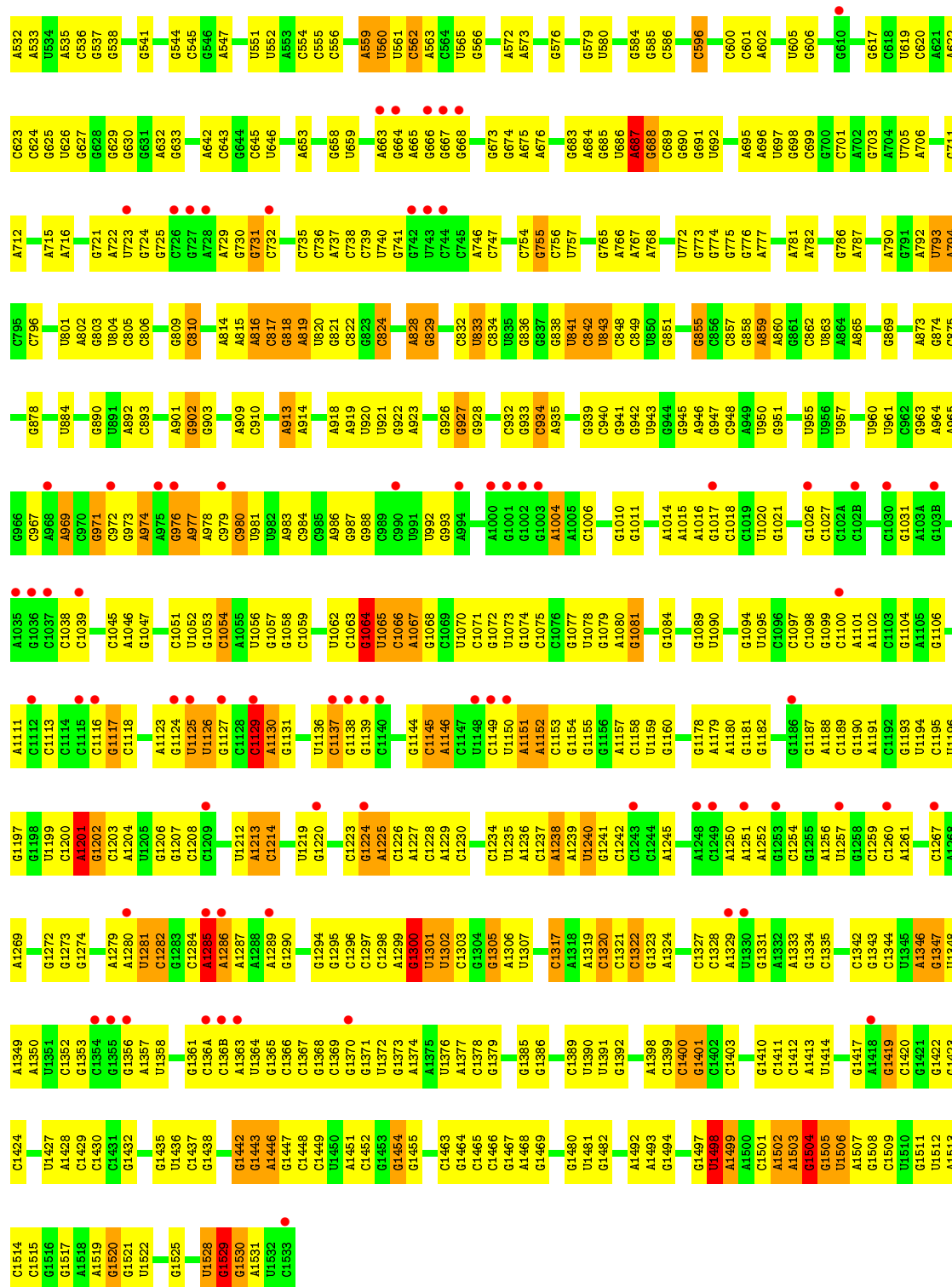
- Molecule 1: ribosomal RNA 16S





• Molecule 1: ribosomal RNA 16S

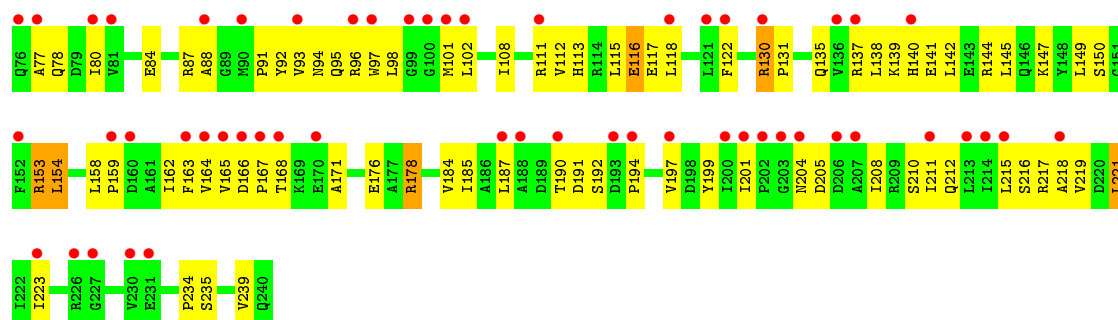




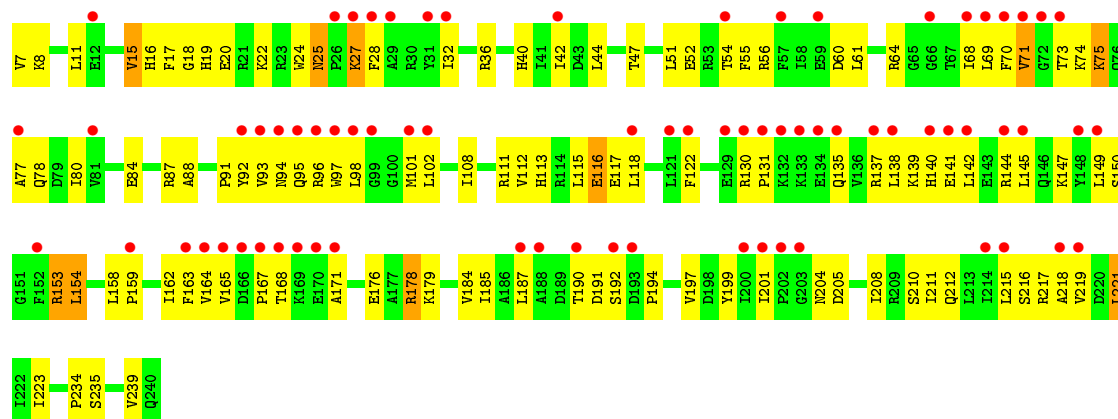
• Molecule 2: 30S ribosomal protein S2

Chain AB:

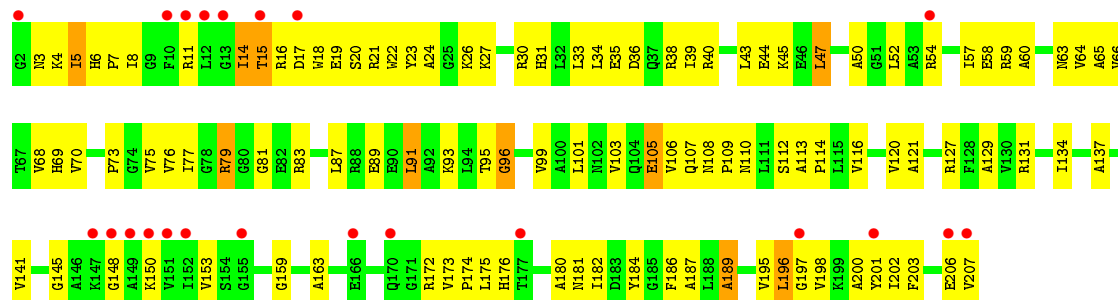




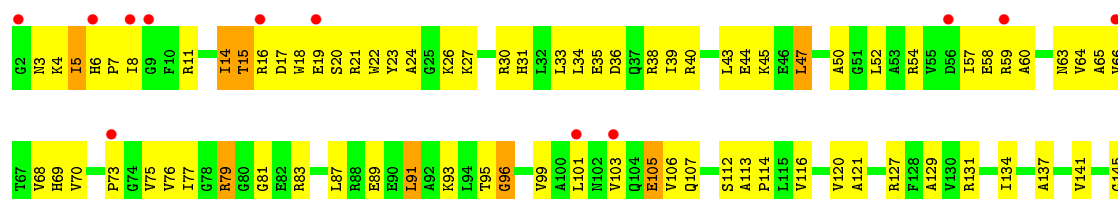
• Molecule 2: 30S ribosomal protein S2

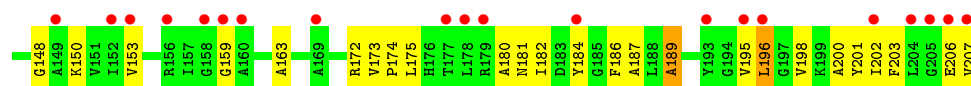


• Molecule 3: 30S ribosomal protein S3

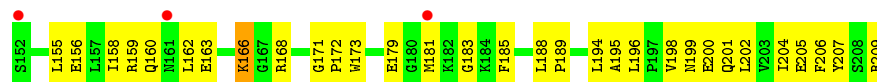


• Molecule 3: 30S ribosomal protein S3

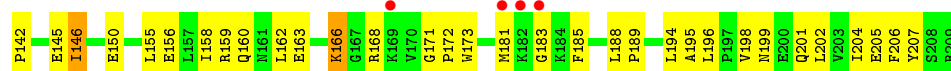
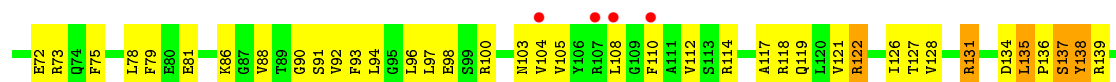
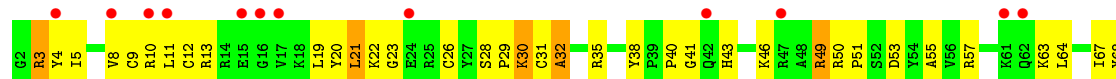




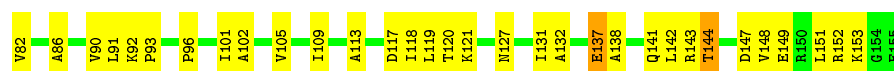
• Molecule 4: 30S ribosomal protein S4



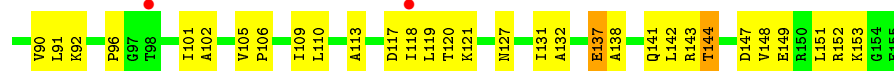
• Molecule 4: 30S ribosomal protein S4



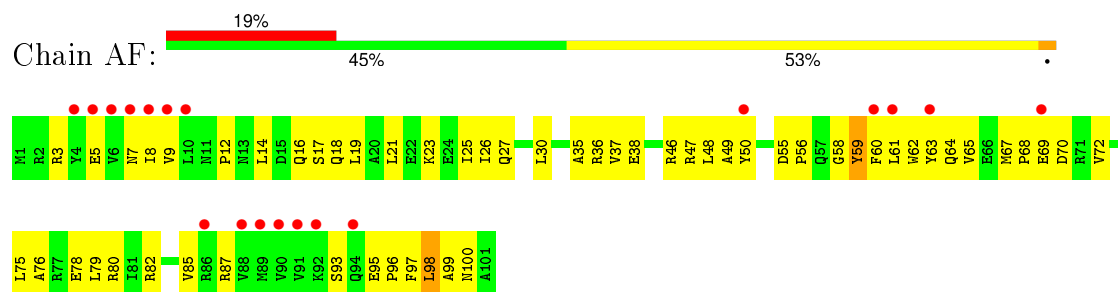
• Molecule 5: 30S ribosomal protein S5



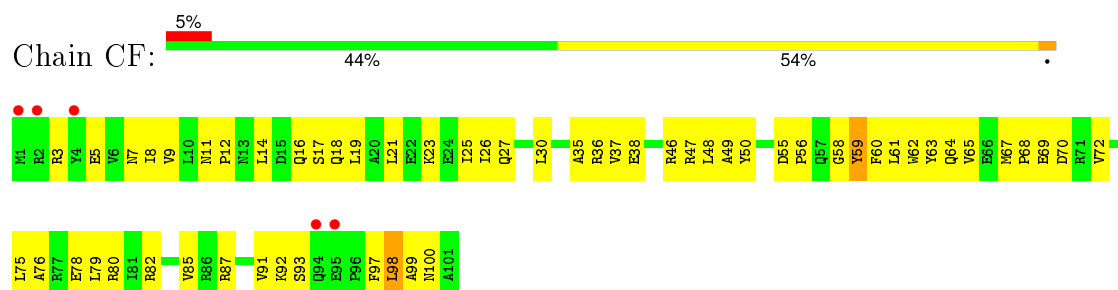
• Molecule 5: 30S ribosomal protein S5



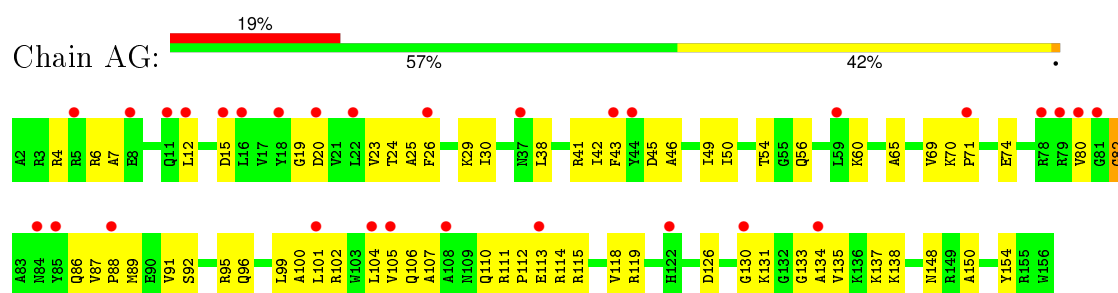
- Molecule 6: 30S ribosomal protein S6



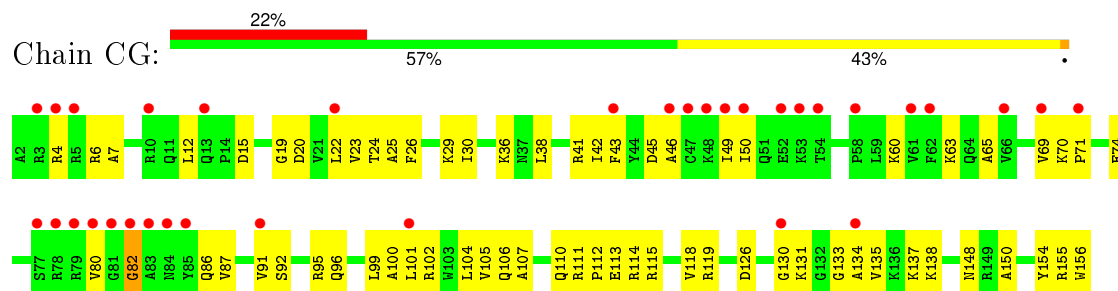
- Molecule 6: 30S ribosomal protein S6



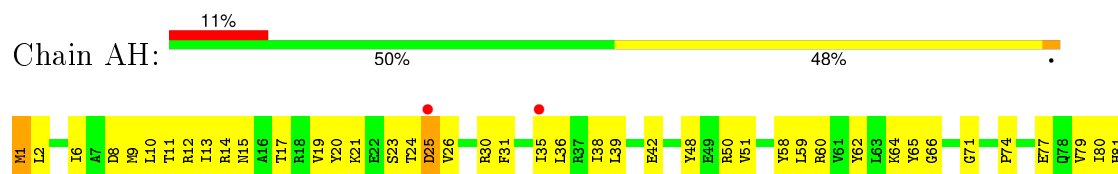
- Molecule 7: 30S ribosomal protein S7

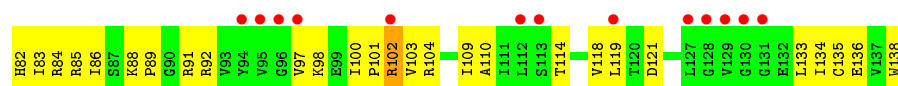


- Molecule 7: 30S ribosomal protein S7

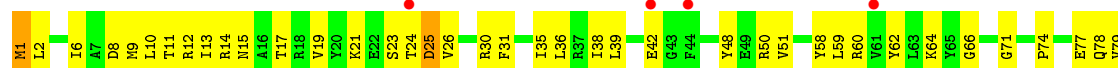


- Molecule 8: 30S ribosomal protein S8

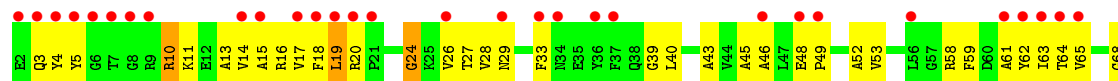
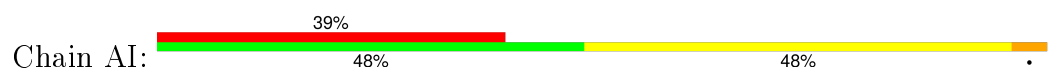




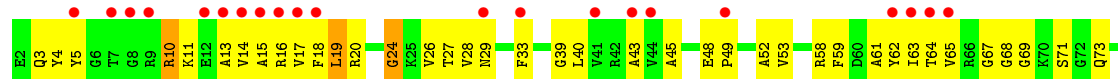
• Molecule 8: 30S ribosomal protein S8



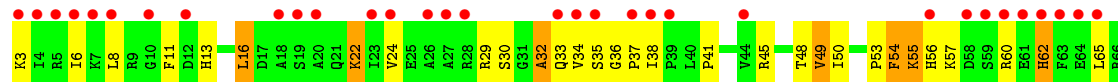
• Molecule 9: 30S ribosomal protein S9



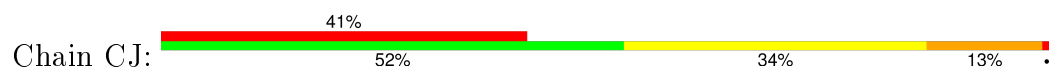
• Molecule 9: 30S ribosomal protein S9

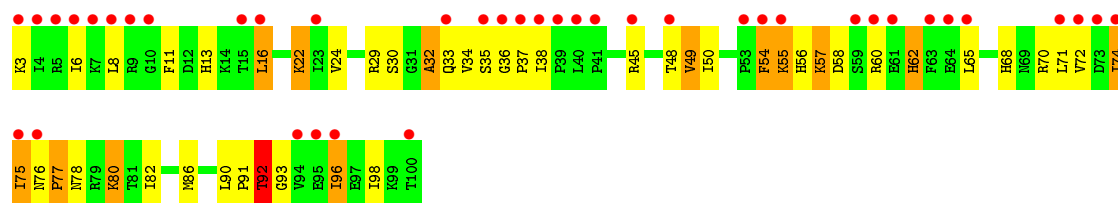


• Molecule 10: 30S ribosomal protein S10

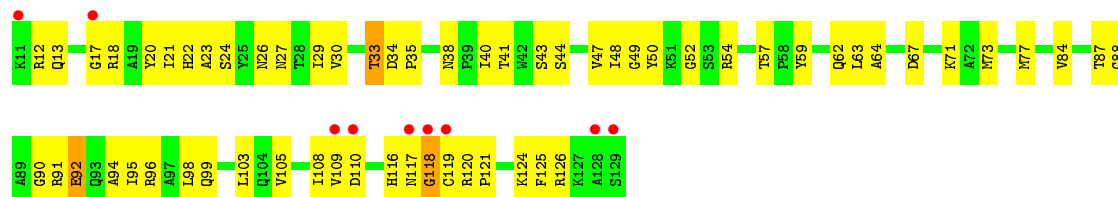


• Molecule 10: 30S ribosomal protein S10

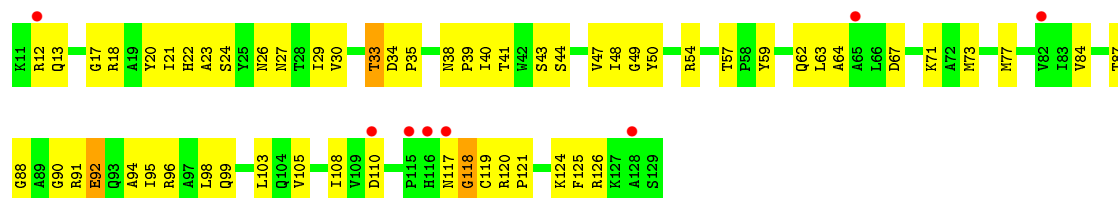




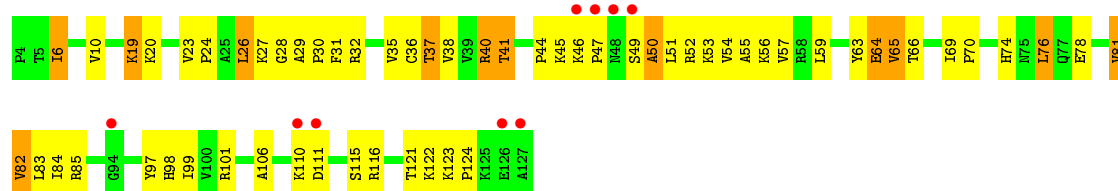
• Molecule 11: 30S ribosomal protein S11



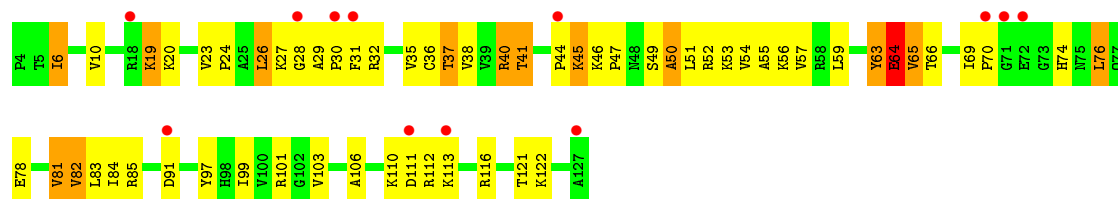
• Molecule 11: 30S ribosomal protein S11



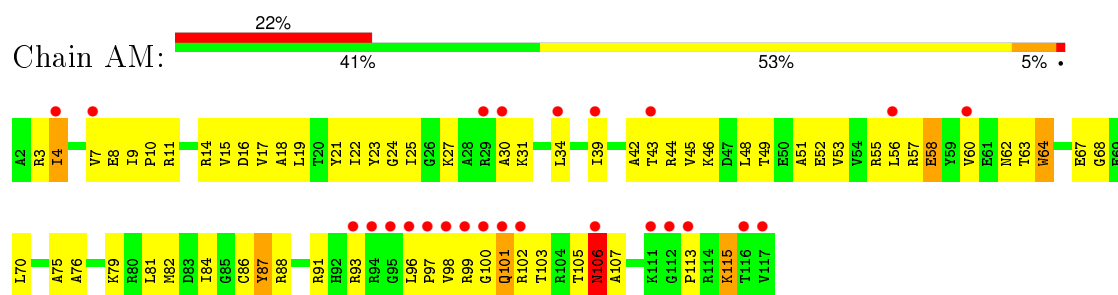
• Molecule 12: 30S ribosomal protein S12



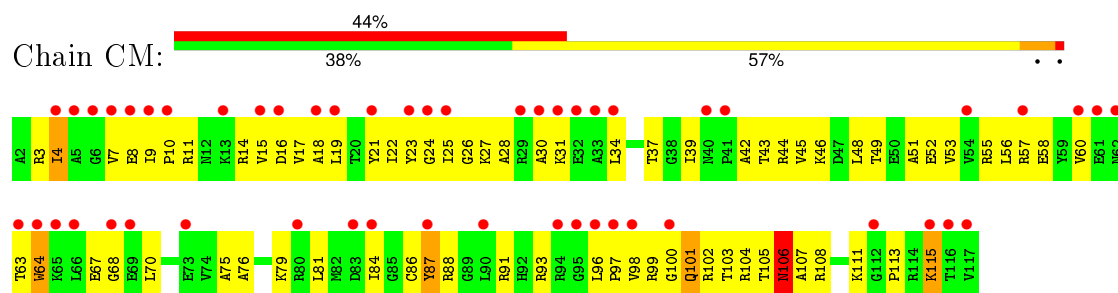
• Molecule 12: 30S ribosomal protein S12



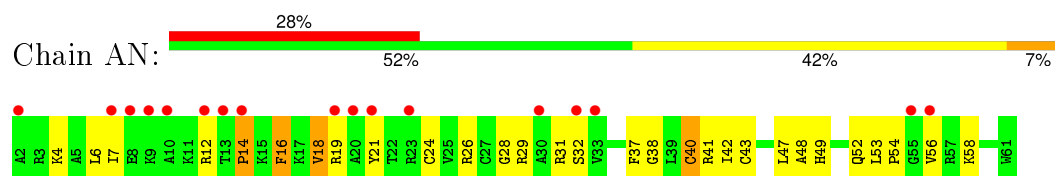
• Molecule 13: 30S ribosomal protein S13



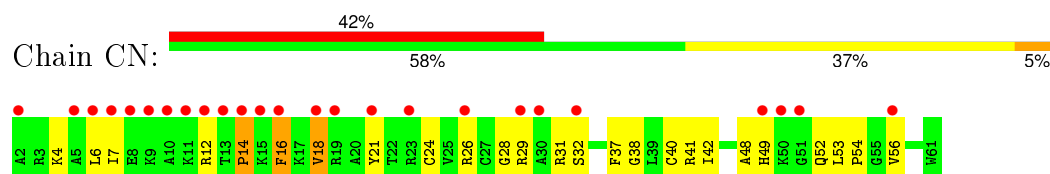
- Molecule 13: 30S ribosomal protein S13



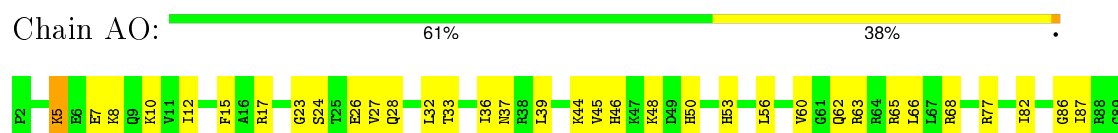
- Molecule 14: 30S ribosomal protein S14 type Z



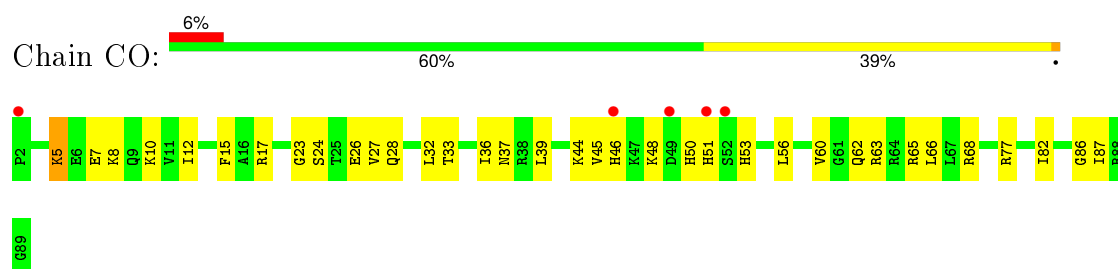
- Molecule 14: 30S ribosomal protein S14 type Z



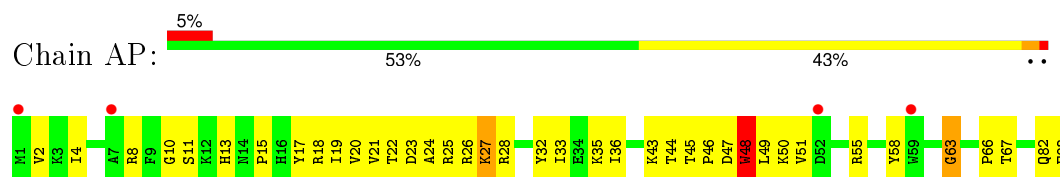
- Molecule 15: 30S ribosomal protein S15



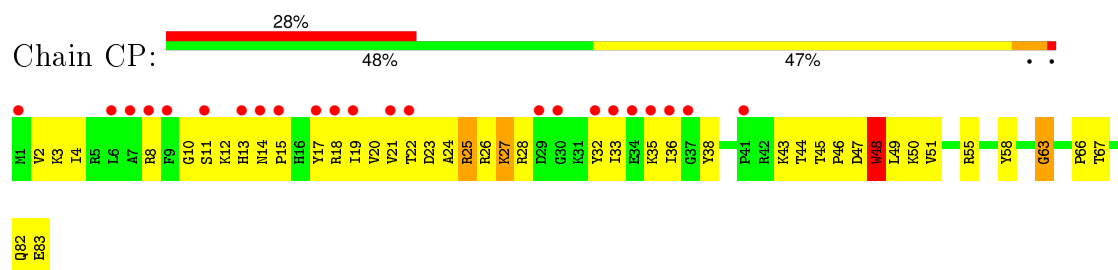
- Molecule 15: 30S ribosomal protein S15



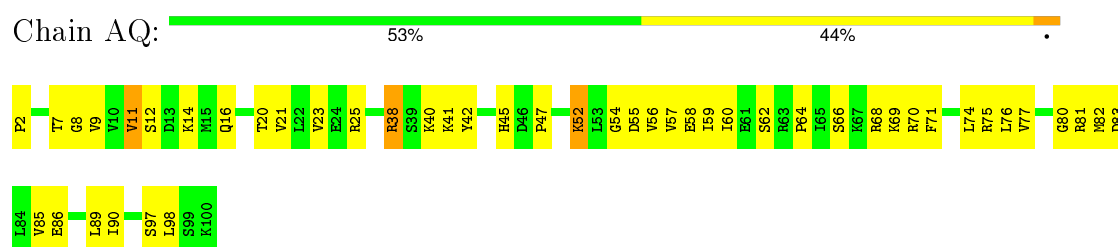
- Molecule 16: 30S ribosomal protein S16



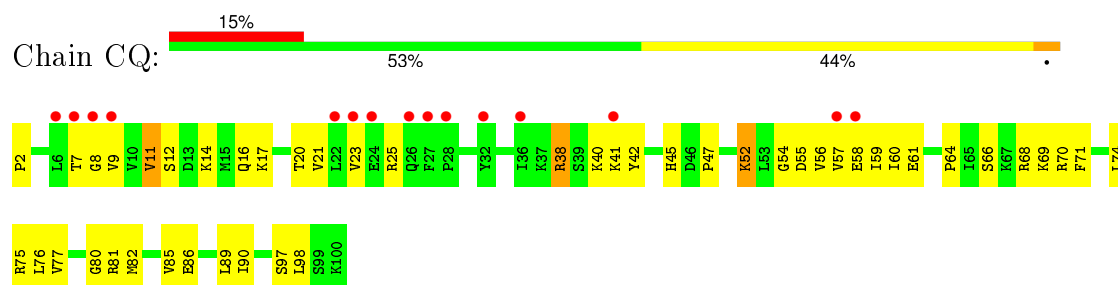
- Molecule 16: 30S ribosomal protein S16



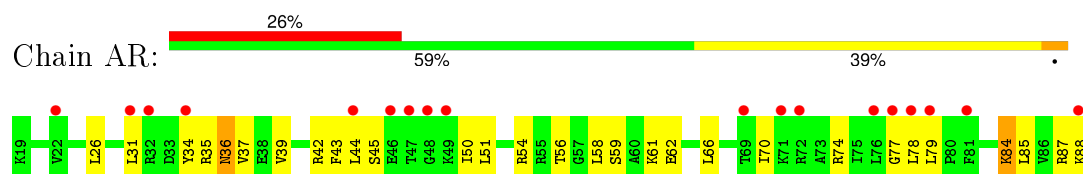
- Molecule 17: 30S ribosomal protein S17



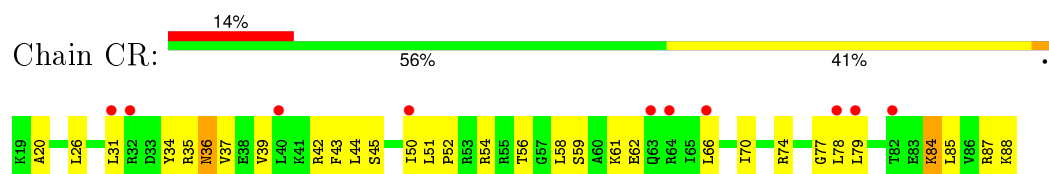
- Molecule 17: 30S ribosomal protein S17



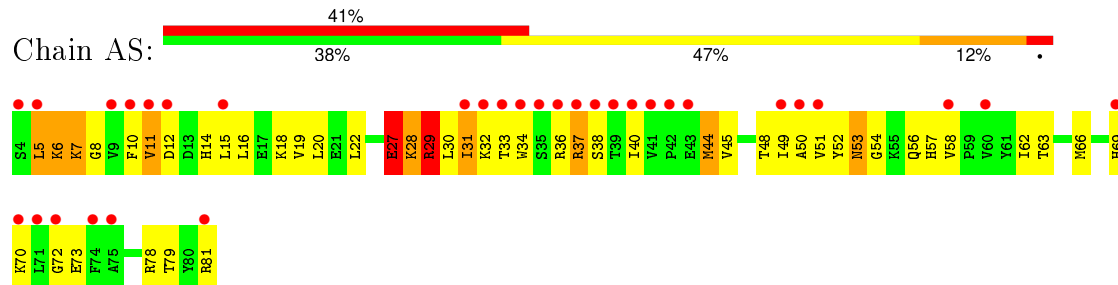
- Molecule 18: 30S ribosomal protein S18



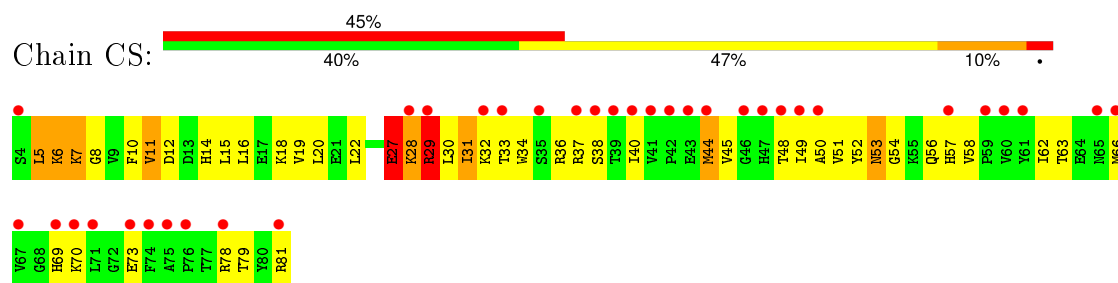
- Molecule 18: 30S ribosomal protein S18



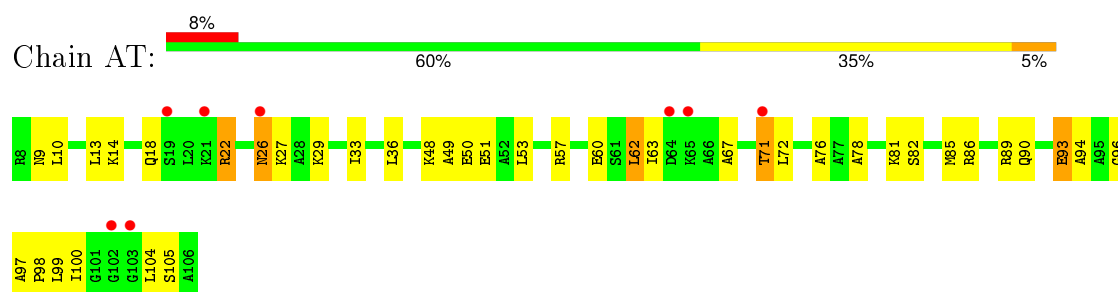
- Molecule 19: 30S ribosomal protein S19



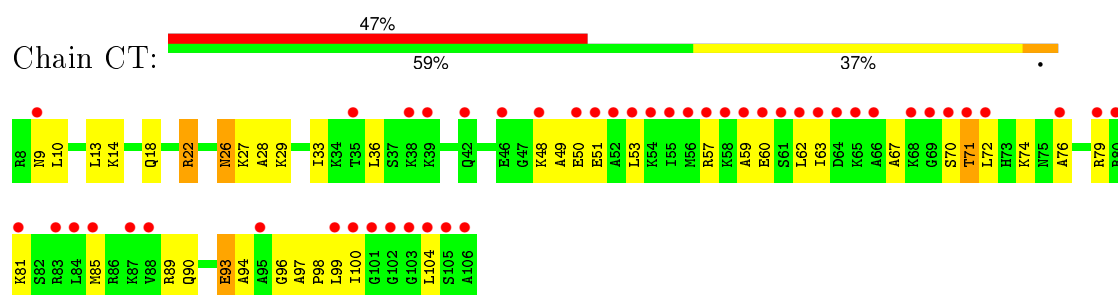
- Molecule 19: 30S ribosomal protein S19



- Molecule 20: 30S ribosomal protein S20



- Molecule 20: 30S ribosomal protein S20

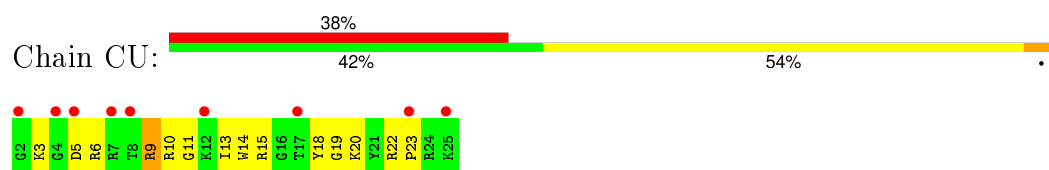


- Molecule 21: 30S ribosomal protein Thx

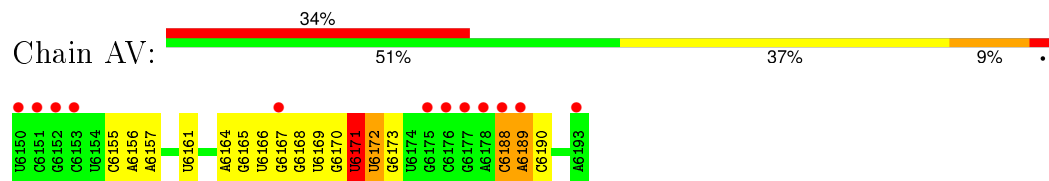


- Molecule 21: 30S ribosomal protein Thx

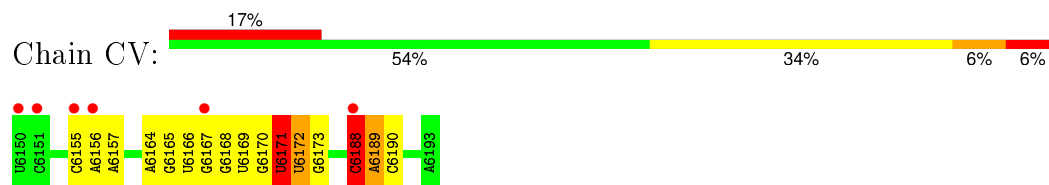




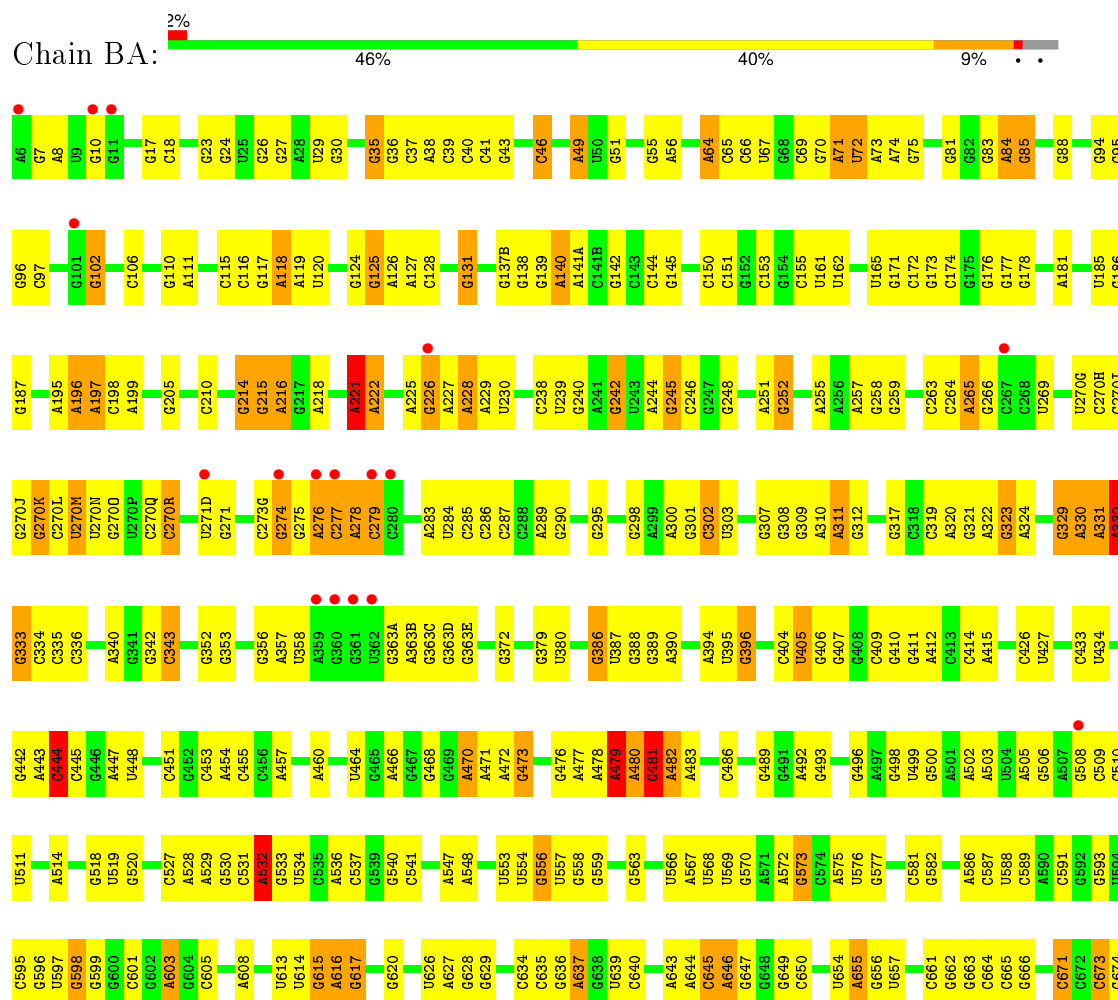
• Molecule 22: domain 3 of PSIC IGR IRES RNA



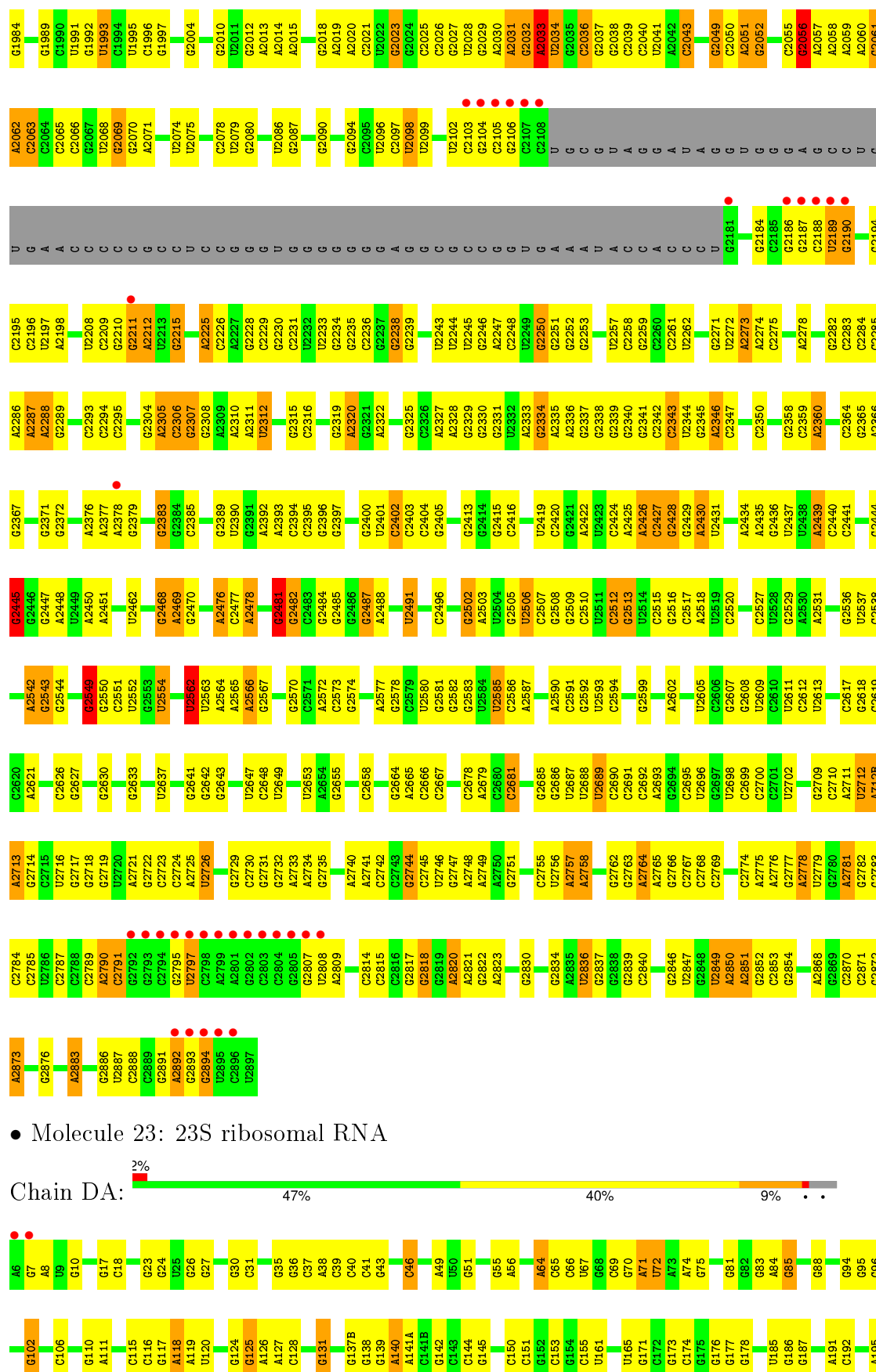
• Molecule 22: domain 3 of PSIC IGR IRES RNA



• Molecule 23: 23S ribosomal RNA

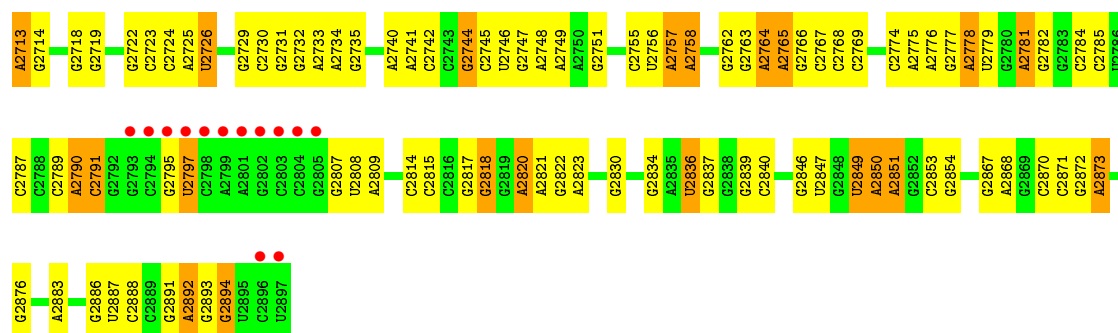


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C1887	U1794	C1684	G1594	U1503	G1413	G1337	G1260	A1155	U	G1002	U930	C939	U773	
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A1900	C1799	U1692	C1599	A1508	G1417	G1343	A1254	G1162	U		G934		A777	G686
A1901	C1800	G1694	C1600	A1509	G1418	G1344	U1255	G1163	A	A1010	G938	C946	G778	
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G1903	A1802	G1696	A1603	A1511	U1420	G1348	C1257	G1165	A	U1012	G940	G848	G780	
	C1804	A1697	C1604		A1427	A1349	G1258	C1166	G	C1013	A941	A849	A781	C692
G1906		A1698		U1516	C1428	C1350			A		G942		A782	
A1913	A1809	A1700	C1607	G1517	G1429	C1351	G1264	G1171	G	U1019	G943	G855	A783	G695
C1914	A1810	A1698	C1608	C1518	U1430	U1352	G1265	G1173	U	A1020	U943	C956	A784	
U1915	G1811	A1699	A1600		U1431	A1353	G1266	A1174	G	A1021	G944	C957	G785	A699
A1916	A1812	C1710	G1611	G1525	C1432	A1354		U1175	C	G1022	A945	U858	G786	G700
U1917	G1813		C1612	G1526	U1433	G1355	C1270	G1176	G	U1023	G946	G859	U787	A706
A1918	G1816	G1726	A1614	A1527	A1434	U1357	G1271	A1177	U	G1024	G947	U860	A788	G707
A1919	A1817	G1727	A1615	U1535	G1435	U1358	A1272	C1178	A	U1025	G948	A861	A789	
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U1930	U1820		A1618	G1538	U1438	G1361	A1275	U1188	G	A1028	A953	A863	G792	U714
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				A1579	A1486		G1326	G1137		A		C914	U827	
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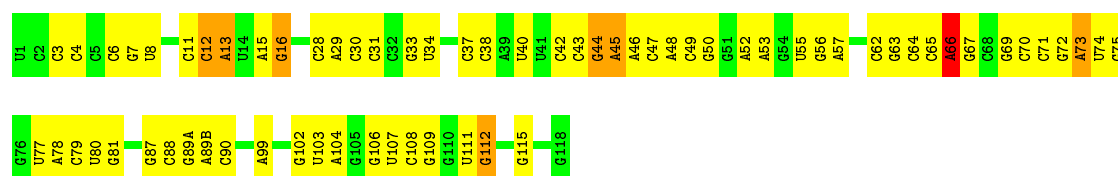
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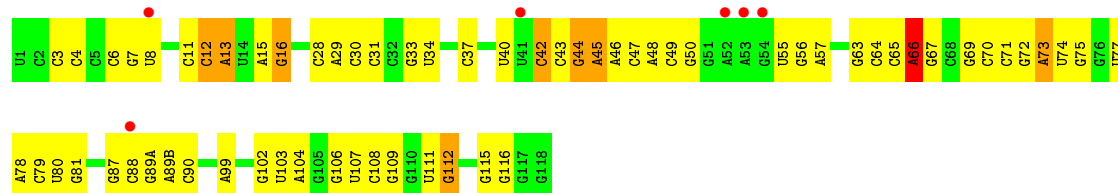
• Molecule 24: 5S ribosomal RNA

Chain BB: 44% 50% 6%



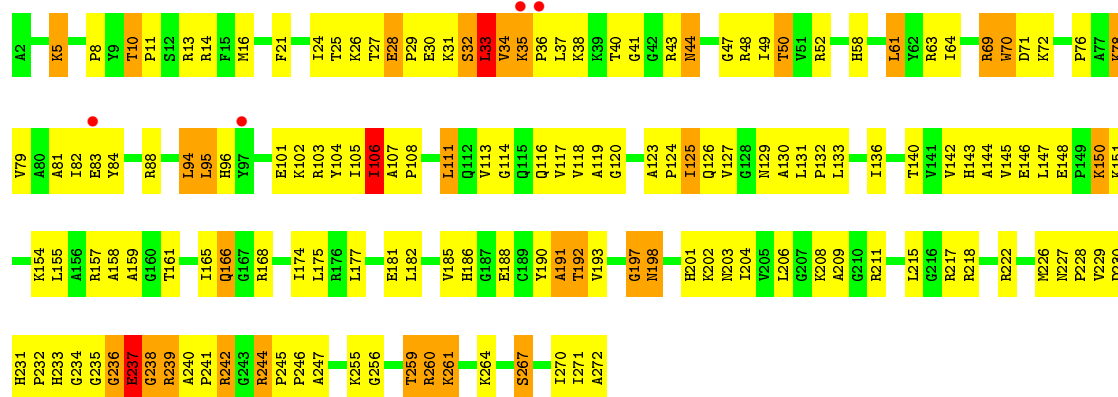
• Molecule 24: 5S ribosomal RNA

Chain DB: 5% 46% 46% 7%



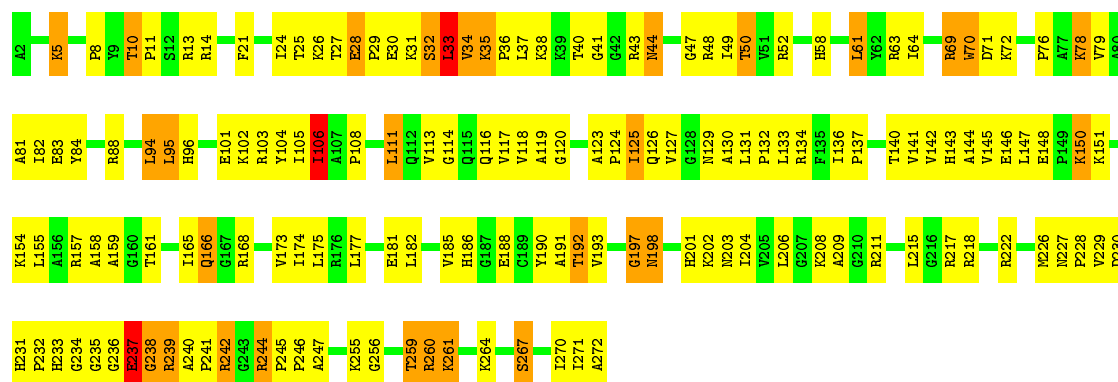
• Molecule 25: 50S ribosomal protein L2

Chain BC: 43% 44% 11%

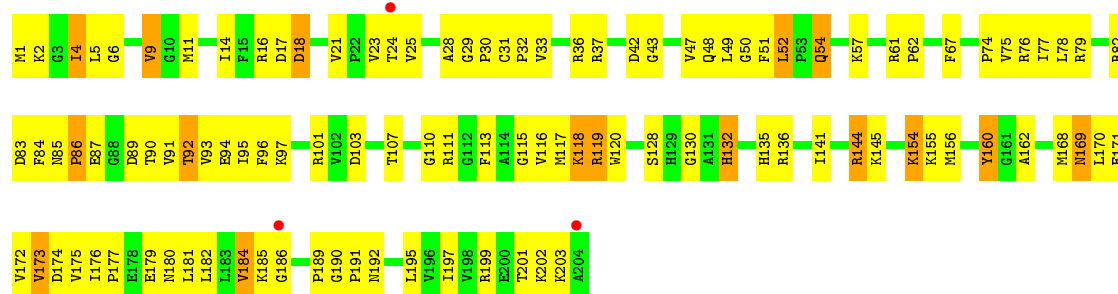


• Molecule 25: 50S ribosomal protein L2

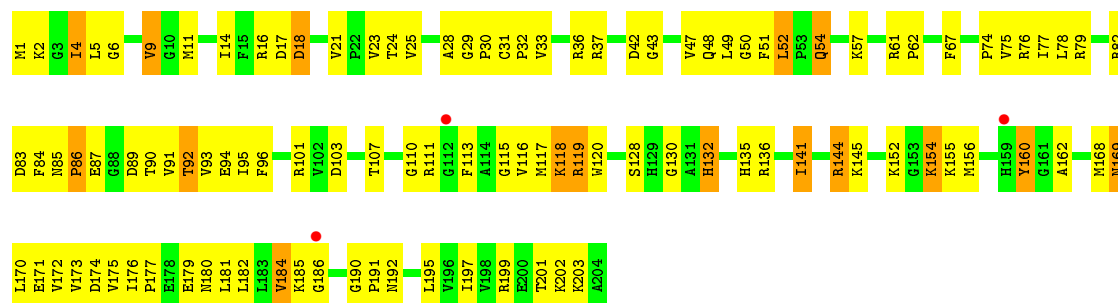
Chain DC: 42% 46% 11%



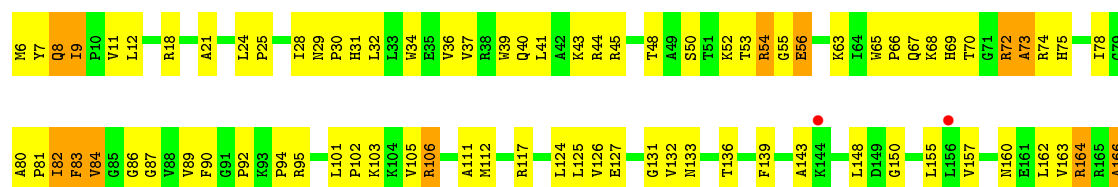
• Molecule 26: 50S ribosomal protein L3

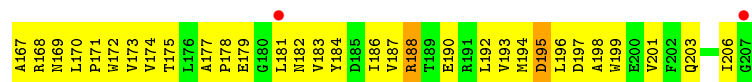


• Molecule 26: 50S ribosomal protein L3

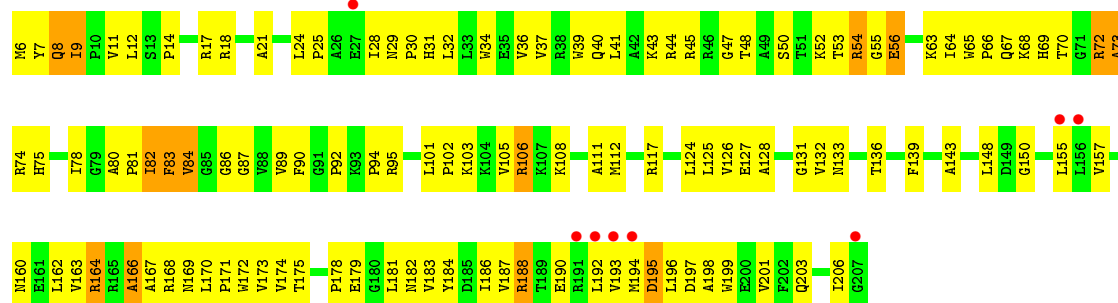
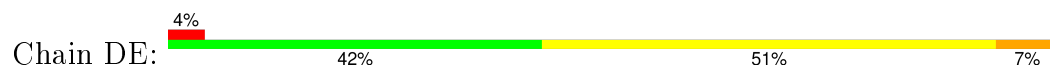


• Molecule 27: 50S ribosomal protein L4

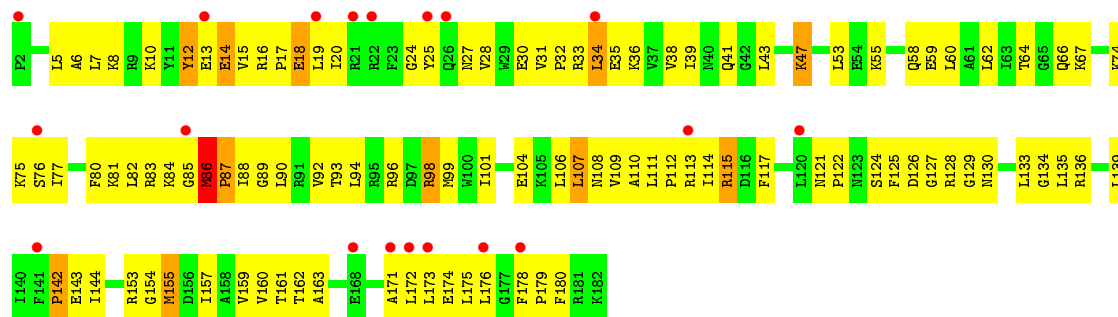




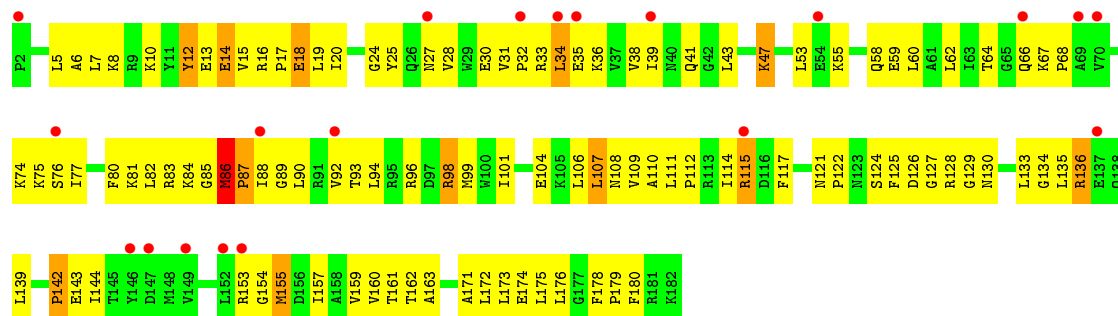
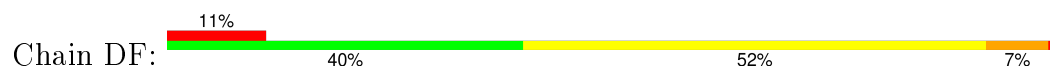
• Molecule 27: 50S ribosomal protein L4



• Molecule 28: 50S ribosomal protein L5



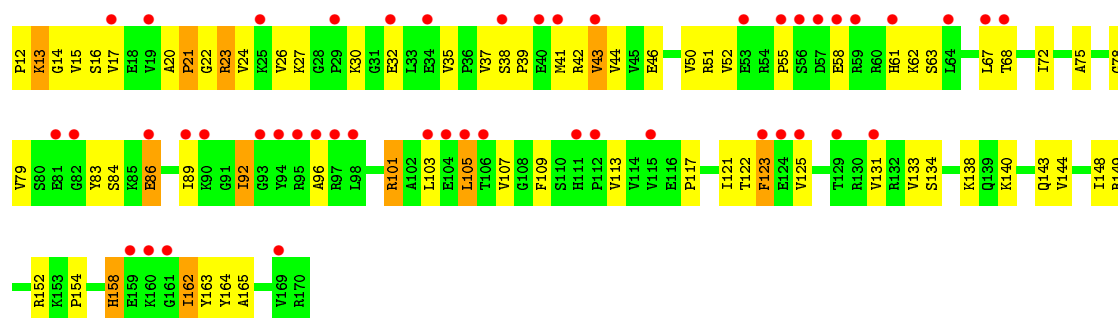
• Molecule 28: 50S ribosomal protein L5



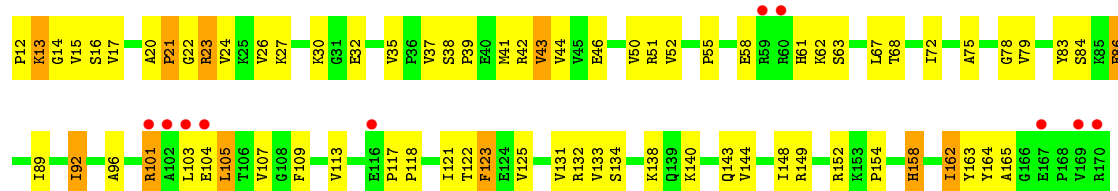
• Molecule 29: 50S ribosomal protein L6



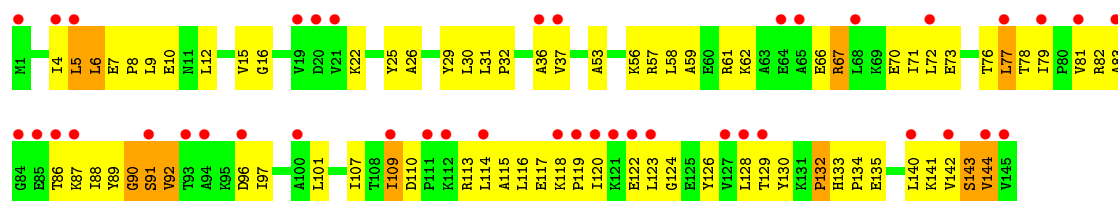




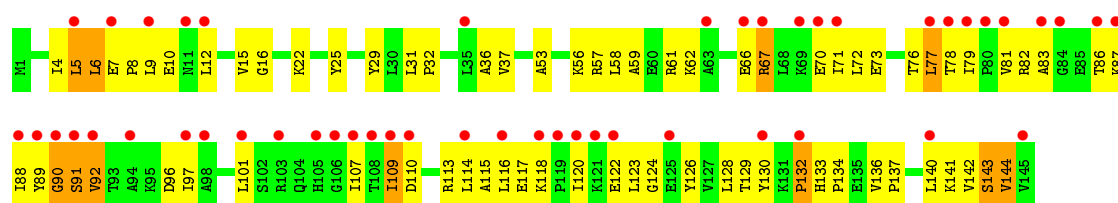
- Molecule 29: 50S ribosomal protein L6



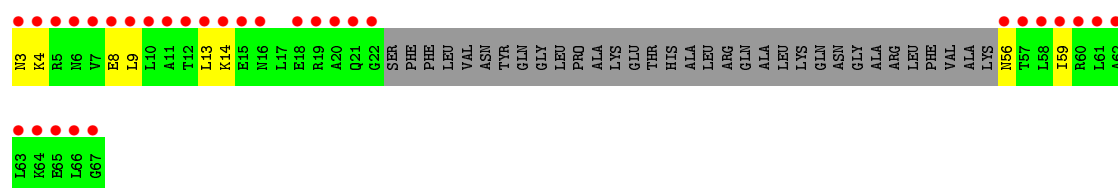
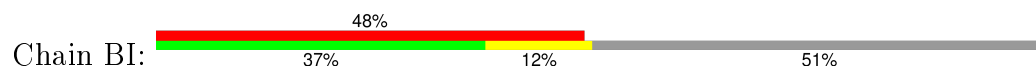
- Molecule 30: 50S ribosomal protein L9



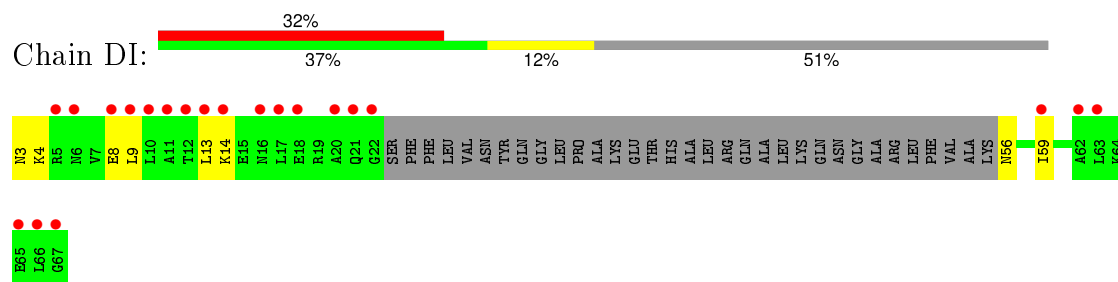
- Molecule 30: 50S ribosomal protein L9



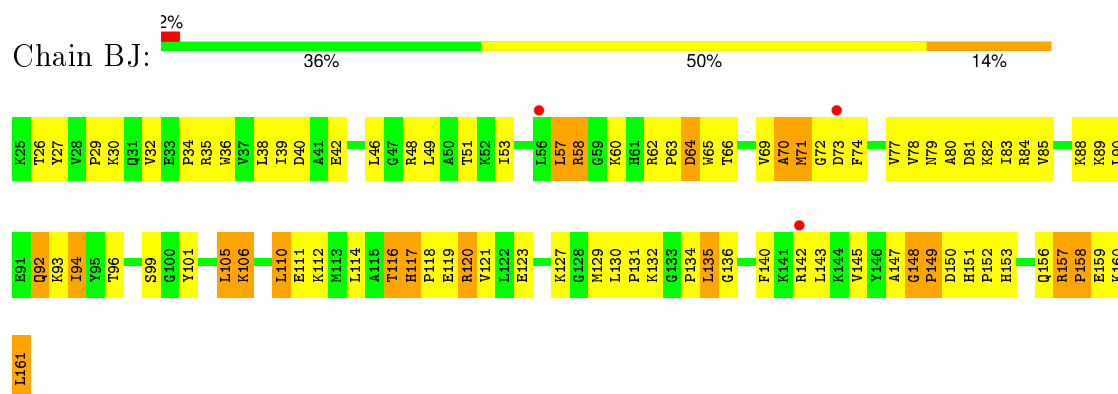
- Molecule 31: 50S ribosomal protein L10



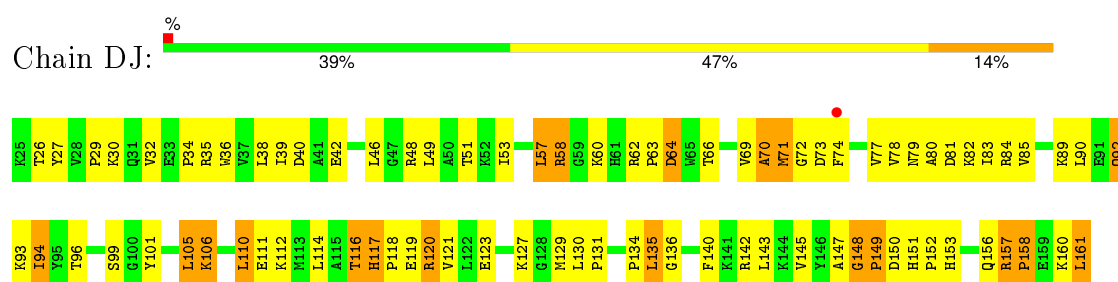
- Molecule 31: 50S ribosomal protein L10



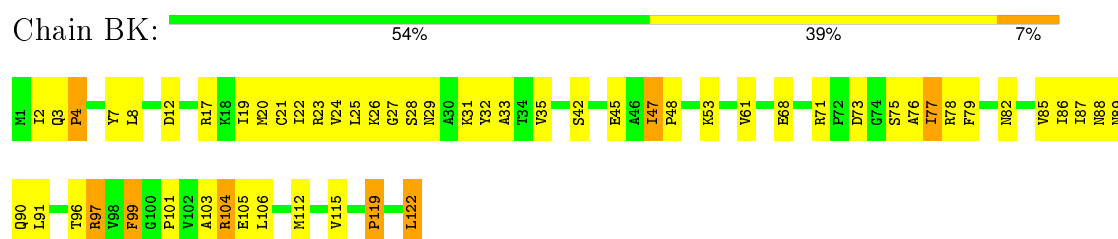
- Molecule 32: 50S ribosomal protein L13



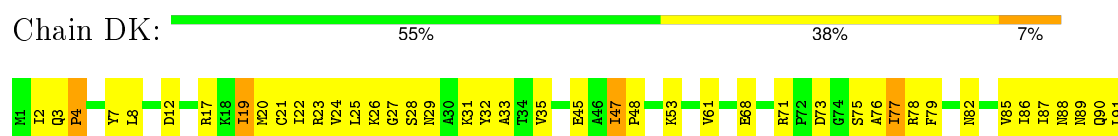
- Molecule 32: 50S ribosomal protein L13

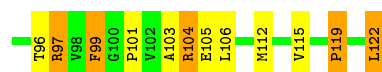


- Molecule 33: 50S ribosomal protein L14

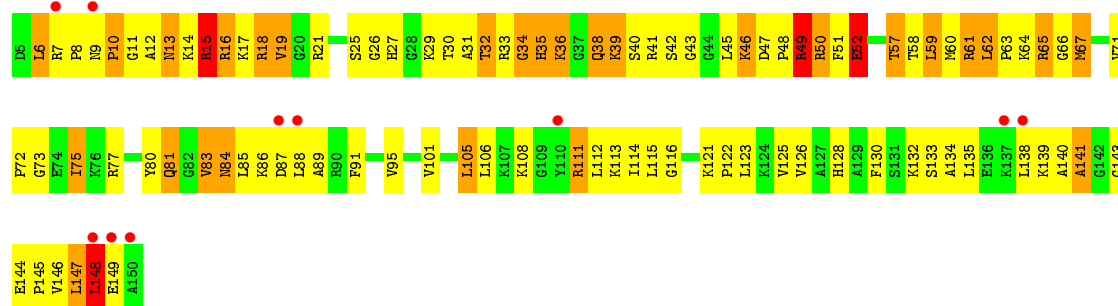


- Molecule 33: 50S ribosomal protein L14

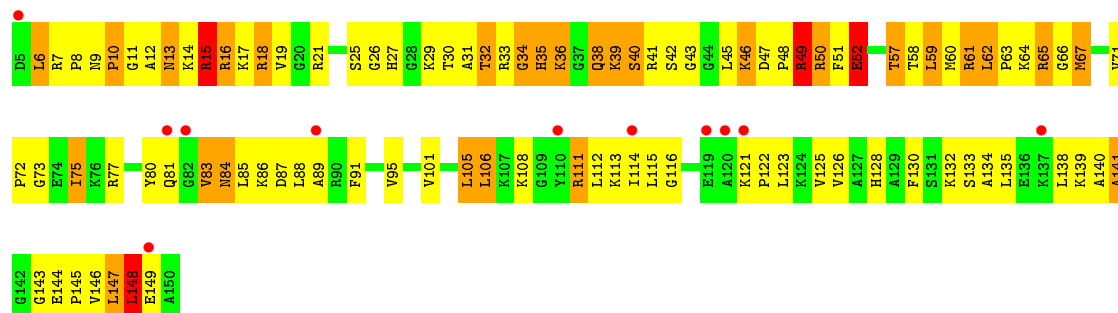




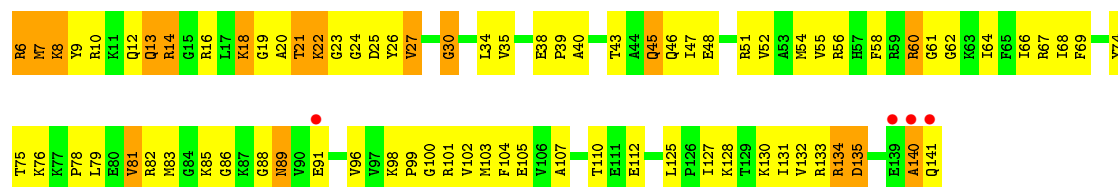
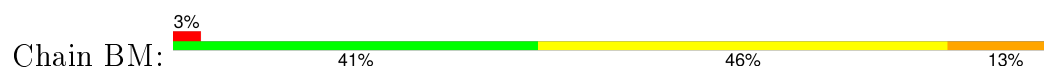
• Molecule 34: 50S ribosomal protein L15



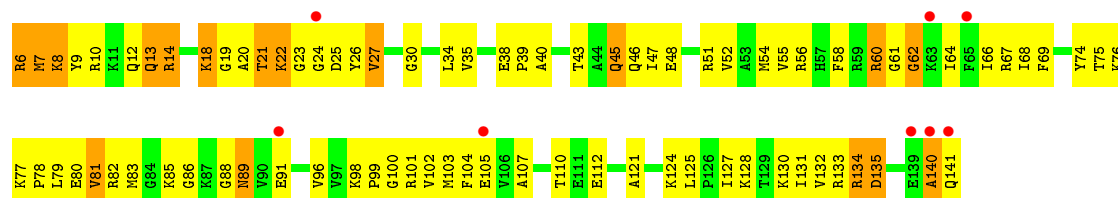
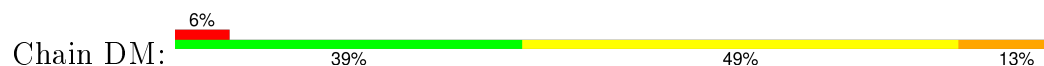
• Molecule 34: 50S ribosomal protein L15

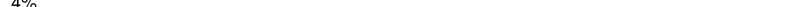


• Molecule 35: 50S ribosomal protein L16



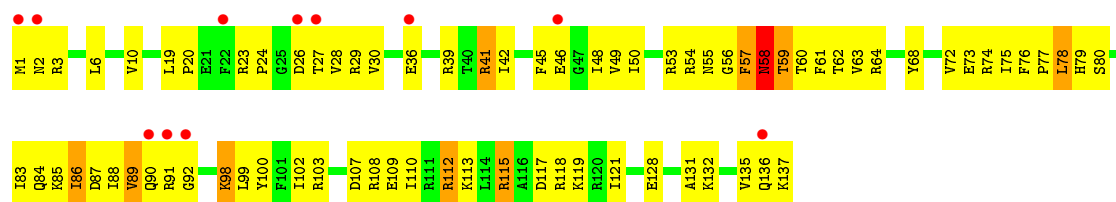
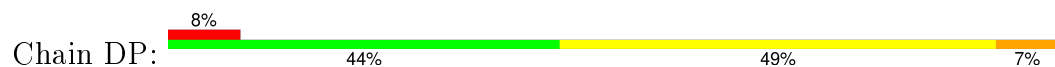
• Molecule 35: 50S ribosomal protein L16



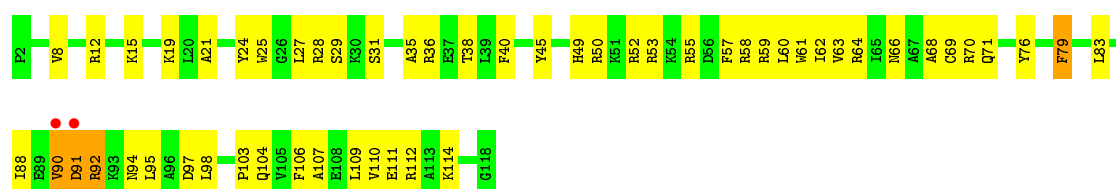
- Chain BP: 



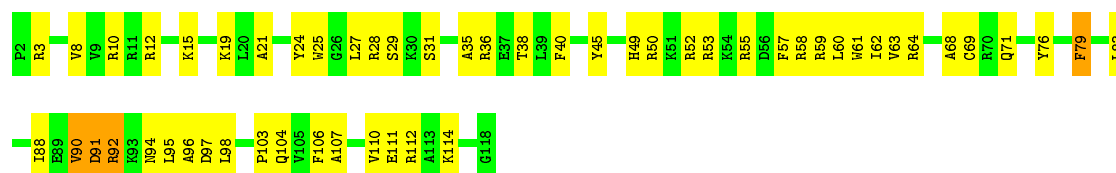
- Molecule 38: 50S ribosomal protein L19



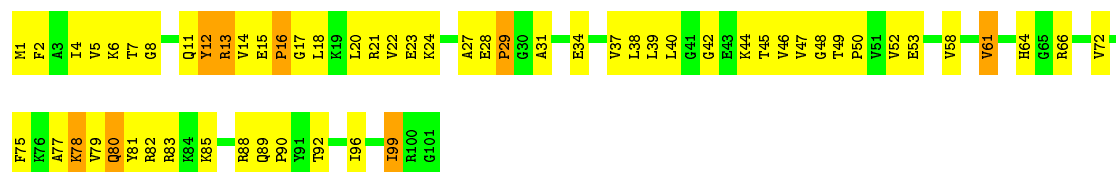
- Molecule 39: 50S ribosomal protein L20



- Molecule 39: 50S ribosomal protein L20



- Molecule 40: 50S ribosomal protein L21

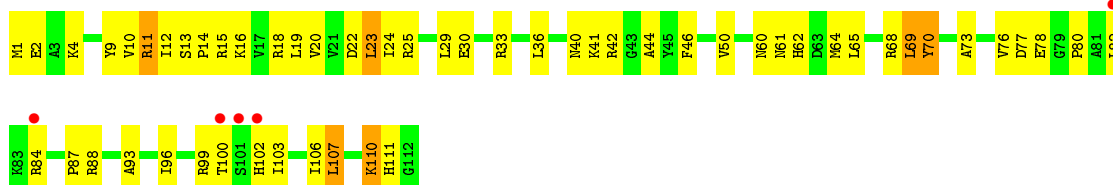


- Molecule 40: 50S ribosomal protein L21

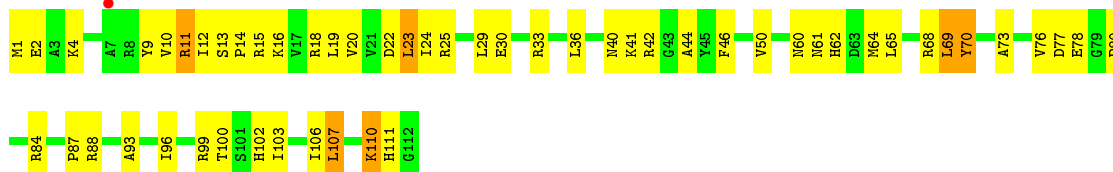




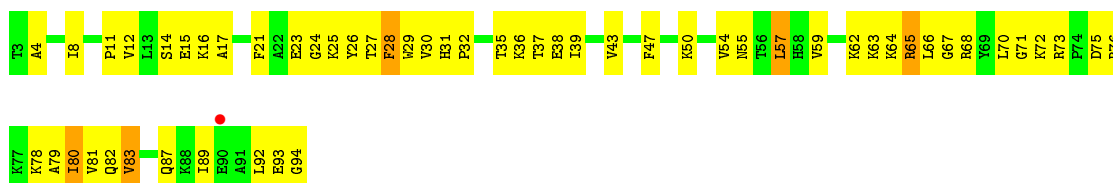
- Molecule 41: 50S ribosomal protein L22



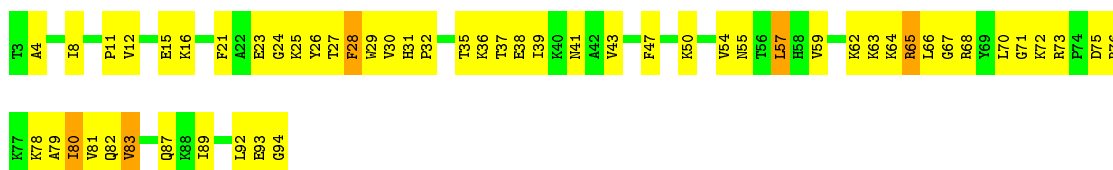
- Molecule 41: 50S ribosomal protein L22



- Molecule 42: 50S ribosomal protein L23

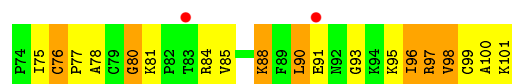


- Molecule 42: 50S ribosomal protein L23

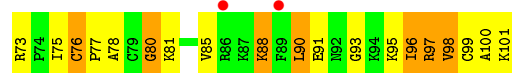
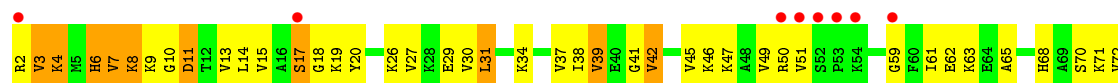


- Molecule 43: 50S ribosomal protein L24

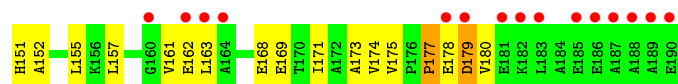
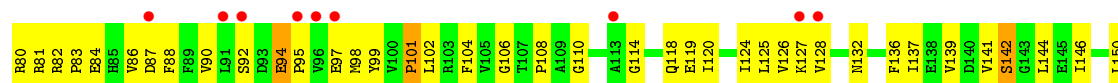
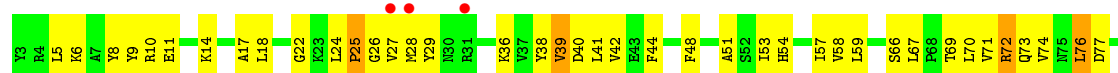




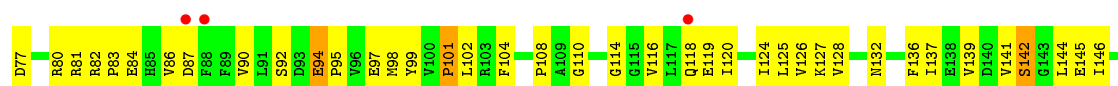
- Molecule 43: 50S ribosomal protein L24



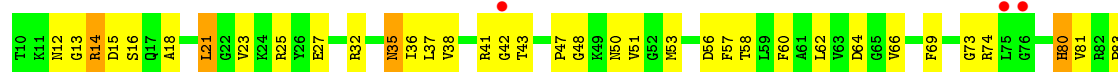
- Molecule 44: 50S ribosomal protein L25



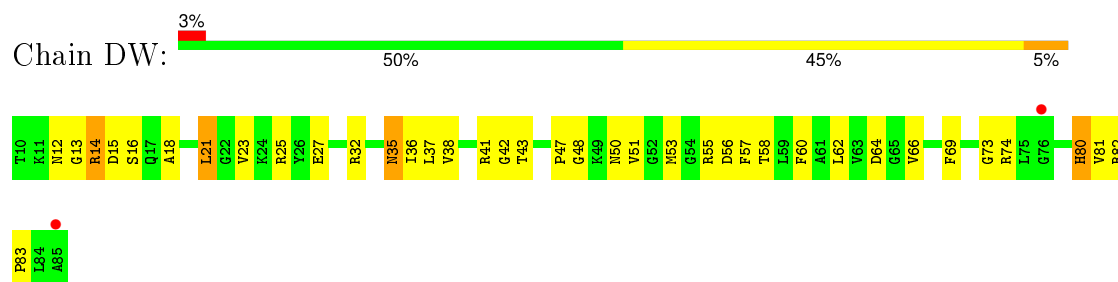
- Molecule 44: 50S ribosomal protein L25



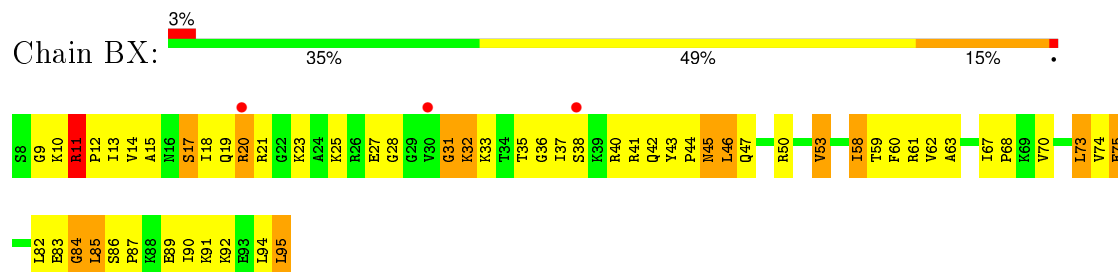
- Molecule 45: 50S ribosomal protein L27



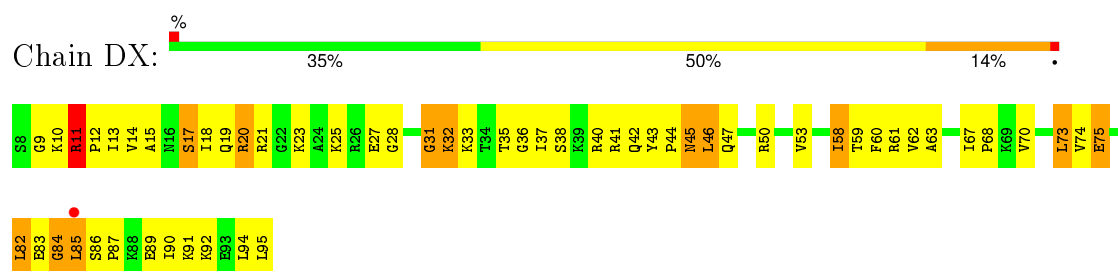
- Molecule 45: 50S ribosomal protein L27



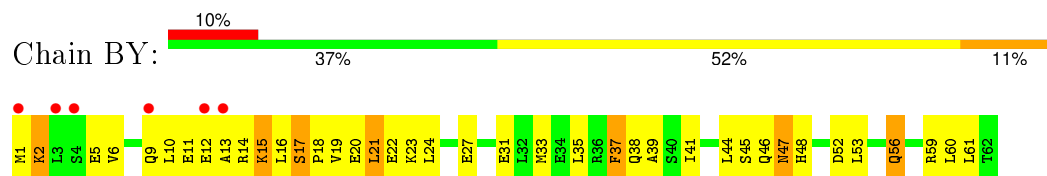
- Molecule 46: 50S ribosomal protein L28



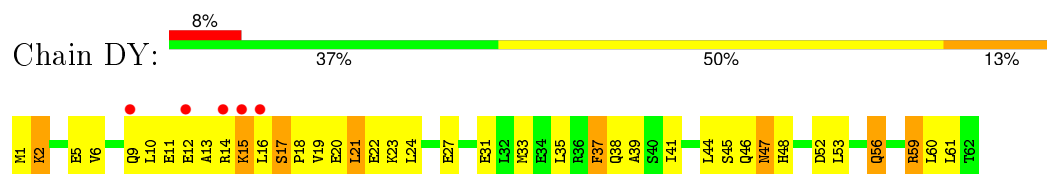
- Molecule 46: 50S ribosomal protein L28



- Molecule 47: 50S ribosomal protein L29



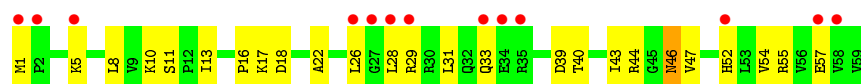
- Molecule 47: 50S ribosomal protein L29



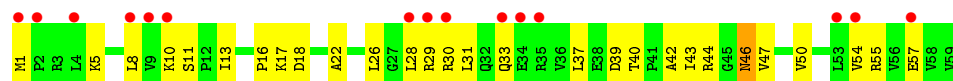
- Molecule 48: 50S ribosomal protein L30



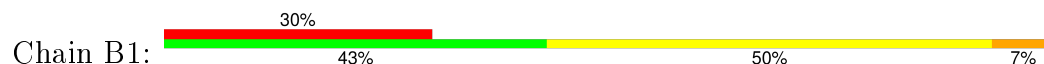




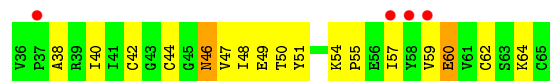
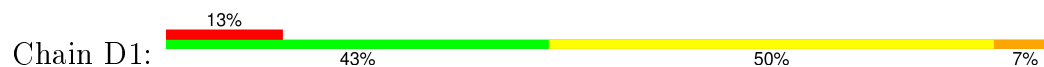
- Molecule 48: 50S ribosomal protein L30



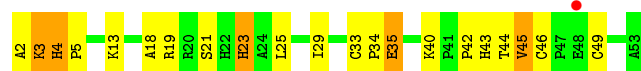
- Molecule 49: 50S ribosomal protein L31



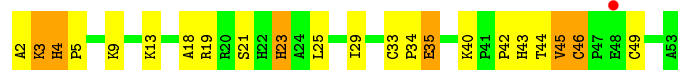
- Molecule 49: 50S ribosomal protein L31



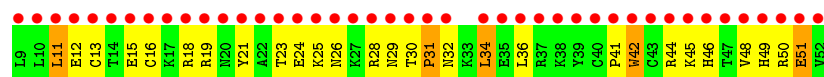
- Molecule 50: 50S ribosomal protein L32



- Molecule 50: 50S ribosomal protein L32

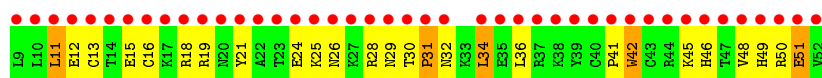


- Molecule 51: 50S ribosomal protein L33

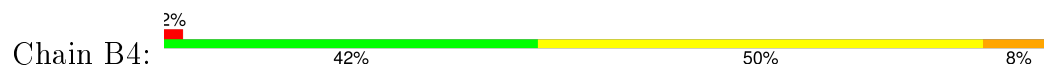


- Molecule 51: 50S ribosomal protein L33





- Molecule 52: 50S ribosomal protein L34



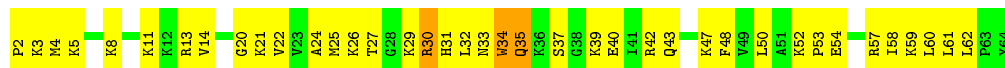
- Molecule 52: 50S ribosomal protein L34



- Molecule 53: 50S ribosomal protein L35



- Molecule 53: 50S ribosomal protein L35



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	211.94Å 455.59Å 618.02Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 3.50 60.01 – 3.49	Depositor EDS
% Data completeness (in resolution range)	99.9 (50.00-3.50) 99.6 (60.01-3.49)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.61 (at 3.49Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.5_2)	Depositor
R, $R_{free}$	0.233 , 0.264 0.229 , 0.262	Depositor DCC
$R_{free}$ test set	7363 reflections (0.99%)	DCC
Wilson B-factor (Å <sup>2</sup> )	106.2	Xtriage
Anisotropy	0.163	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.25 , 109.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.38$ , $\langle L^2 \rangle = 0.21$	Xtriage
Outliers	1 of 746568 reflections (0.000%)	Xtriage
$F_o, F_c$ correlation	0.91	EDS
Total number of atoms	283641	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	127.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.63% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: ZN, MG

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	AA	0.45	0/36238	0.90	34/56561 (0.1%)
1	CA	0.44	0/36238	0.90	34/56561 (0.1%)
2	AB	0.21	0/1936	0.38	0/2609
2	CB	0.21	0/1936	0.38	0/2609
3	AC	0.21	0/1637	0.37	0/2205
3	CC	0.21	0/1637	0.37	0/2205
4	AD	0.24	0/1733	0.40	0/2318
4	CD	0.23	0/1733	0.39	0/2318
5	AE	0.24	0/1172	0.41	0/1576
5	CE	0.23	0/1172	0.40	0/1576
6	AF	0.23	0/856	0.42	0/1154
6	CF	0.23	0/856	0.43	0/1154
7	AG	0.21	0/1276	0.36	0/1709
7	CG	0.22	0/1276	0.36	0/1709
8	AH	0.23	0/1136	0.41	0/1527
8	CH	0.22	0/1136	0.41	0/1527
9	AI	0.22	0/1029	0.38	0/1378
9	CI	0.22	0/1029	0.38	0/1378
10	AJ	0.21	0/808	0.39	0/1085
10	CJ	0.21	0/808	0.39	0/1085
11	AK	0.23	0/900	0.40	0/1213
11	CK	0.23	0/900	0.40	0/1213
12	AL	0.25	0/987	0.46	0/1320
12	CL	0.24	0/987	0.46	0/1320
13	AM	0.24	0/939	0.41	0/1258
13	CM	0.24	0/939	0.41	0/1258
14	AN	0.23	0/501	0.37	0/664
14	CN	0.22	0/501	0.37	0/664
15	AO	0.24	0/745	0.38	0/992
15	CO	0.24	0/745	0.38	0/992
16	AP	0.36	1/717 (0.1%)	0.44	0/963
16	CP	0.44	1/717 (0.1%)	0.45	0/963

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z  >5	RMSZ	# Z  >5
17	AQ	0.24	0/837	0.41	0/1117
17	CQ	0.23	0/837	0.40	0/1117
18	AR	0.24	0/579	0.42	0/768
18	CR	0.23	0/579	0.42	0/768
19	AS	0.20	0/643	0.37	0/865
19	CS	0.21	0/643	0.37	0/865
20	AT	0.23	0/764	0.39	0/1006
20	CT	0.22	0/764	0.39	0/1006
21	AU	0.20	0/213	0.37	0/277
21	CU	0.19	0/213	0.37	0/277
22	AV	0.43	0/821	0.86	2/1275 (0.2%)
22	CV	0.44	0/821	0.86	2/1275 (0.2%)
23	BA	0.51	0/66575	1.03	115/103930 (0.1%)
23	DA	0.54	0/66575	1.04	119/103930 (0.1%)
24	BB	0.44	0/2853	0.92	1/4451 (0.0%)
24	DB	0.44	0/2853	0.93	2/4451 (0.0%)
25	BC	0.33	0/2155	0.51	0/2905
25	DC	0.34	0/2155	0.52	0/2905
26	BD	0.27	0/1597	0.48	0/2153
26	DD	0.27	0/1597	0.48	0/2153
27	BE	0.29	0/1622	0.46	0/2194
27	DE	0.31	0/1622	0.47	0/2194
28	BF	0.23	0/1500	0.42	0/2017
28	DF	0.23	0/1500	0.42	0/2017
29	BG	0.22	0/1246	0.42	0/1682
29	DG	0.24	0/1246	0.43	0/1682
30	BH	0.29	0/1148	0.46	0/1552
30	DH	0.31	0/1148	0.47	0/1552
31	BI	0.21	0/252	0.38	0/333
31	DI	0.22	0/252	0.38	0/333
32	BJ	0.26	0/1124	0.47	0/1515
32	DJ	0.27	0/1124	0.47	0/1515
33	BK	0.27	0/942	0.48	0/1268
33	DK	0.28	0/942	0.49	0/1268
34	BL	0.30	0/1131	0.56	0/1504
34	DL	0.32	0/1131	0.57	0/1504
35	BM	0.30	0/1099	0.49	0/1468
35	DM	0.30	0/1099	0.50	0/1468
36	BN	0.26	0/974	0.45	0/1302
36	DN	0.27	0/974	0.45	0/1302
37	BO	0.23	0/779	0.42	0/1036
37	DO	0.24	0/779	0.42	0/1036
38	BP	0.27	0/1158	0.44	0/1544

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z  >5	RMSZ	# Z  >5
38	DP	0.27	0/1158	0.44	0/1544
39	BQ	0.29	0/970	0.46	0/1290
39	DQ	0.31	0/970	0.47	0/1290
40	BR	0.27	0/790	0.45	0/1057
40	DR	0.29	0/790	0.46	0/1057
41	BS	0.31	0/902	0.51	0/1209
41	DS	0.30	0/902	0.50	0/1209
42	BT	0.30	0/740	0.49	0/993
42	DT	0.33	0/740	0.50	0/993
43	BU	0.25	0/789	0.44	0/1051
43	DU	0.26	0/789	0.45	0/1051
44	BV	0.23	0/1524	0.44	0/2068
44	DV	0.23	0/1524	0.45	0/2068
45	BW	0.26	0/613	0.43	0/816
45	DW	0.27	0/613	0.43	0/816
46	BX	0.30	0/702	0.56	0/932
46	DX	0.31	0/702	0.57	0/932
47	BY	0.29	0/523	0.52	0/690
47	DY	0.31	0/523	0.54	0/690
48	BZ	0.23	0/473	0.41	0/634
48	DZ	0.24	0/473	0.41	0/634
49	B1	0.20	0/229	0.38	0/309
49	D1	0.20	0/229	0.38	0/309
50	B2	0.28	0/419	0.51	0/567
50	D2	0.28	0/419	0.51	0/567
51	B3	0.21	0/388	0.40	0/518
51	D3	0.21	0/388	0.40	0/518
52	B4	0.34	0/427	0.52	0/561
52	D4	0.38	0/427	0.53	0/561
53	B5	0.31	0/516	0.50	0/679
53	D5	0.32	0/516	0.51	0/679
All	All	0.44	2/305254 (0.0%)	0.87	309/456136 (0.1%)

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
28	BF	0	1
28	DF	0	1
34	BL	0	1

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	#Chirality outliers	#Planarity outliers
34	DL	0	1
All	All	0	4

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
16	CP	48	TRP	CG-CD1	7.51	1.47	1.36
16	AP	48	TRP	CG-CD1	6.09	1.45	1.36

All (309) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
23	BA	945	A	C1'-O4'-C4'	-10.48	101.52	109.90
23	BA	1786	A	C1'-O4'-C4'	-10.25	101.70	109.90
23	DA	676	A	C1'-O4'-C4'	-9.79	102.07	109.90
23	DA	1786	A	C1'-O4'-C4'	-9.78	102.07	109.90
23	DA	945	A	C1'-O4'-C4'	-9.73	102.11	109.90
23	BA	676	A	C1'-O4'-C4'	-9.48	102.32	109.90
23	BA	2818	G	C1'-O4'-C4'	-9.24	102.51	109.90
23	DA	945	A	O4'-C1'-N9	9.13	115.50	108.20
23	DA	2818	G	C1'-O4'-C4'	-9.09	102.63	109.90
1	CA	99	C	P-O3'-C3'	9.00	130.50	119.70
23	BA	945	A	O4'-C1'-N9	8.85	115.28	108.20
1	AA	99	C	P-O3'-C3'	8.79	130.25	119.70
23	DA	265	A	C3'-C2'-C1'	-8.72	94.53	101.50
23	BA	265	A	C3'-C2'-C1'	-8.68	94.55	101.50
23	DA	761	A	N1-C6-N6	8.67	123.80	118.60
23	DA	748	G	C1'-O4'-C4'	-8.53	103.08	109.90
23	DA	221	A	P-O3'-C3'	8.12	129.45	119.70
23	DA	1786	A	C3'-C2'-C1'	-8.09	95.03	101.50
23	BA	1786	A	C3'-C2'-C1'	-8.01	95.09	101.50
23	BA	761	A	N1-C6-N6	7.98	123.39	118.60
1	CA	266	G	P-O3'-C3'	7.96	129.26	119.70
1	AA	266	G	P-O3'-C3'	7.94	129.23	119.70
23	BA	221	A	P-O3'-C3'	7.85	129.12	119.70
23	BA	1614	A	C1'-O4'-C4'	-7.84	103.63	109.90
23	BA	481	G	P-O3'-C3'	7.74	128.98	119.70
23	BA	748	G	C1'-O4'-C4'	-7.71	103.73	109.90
23	DA	1614	A	C1'-O4'-C4'	-7.70	103.74	109.90
23	DA	1937	A	P-O3'-C3'	7.66	128.89	119.70
1	AA	1498	U	P-O3'-C3'	7.63	128.86	119.70
23	BA	1937	A	P-O3'-C3'	7.59	128.81	119.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
23	DA	933	A	O4'-C1'-N9	7.51	114.21	108.20
23	DA	481	G	P-O3'-C3'	7.49	128.69	119.70
1	AA	243	A	P-O3'-C3'	7.42	128.61	119.70
23	DA	945	A	C3'-C2'-C1'	-7.42	95.56	101.50
1	CA	1529	G	C1'-O4'-C4'	-7.38	103.99	109.90
23	BA	933	A	O4'-C1'-N9	7.32	114.06	108.20
1	CA	243	A	P-O3'-C3'	7.25	128.40	119.70
1	CA	1498	U	P-O3'-C3'	7.23	128.38	119.70
23	BA	945	A	C3'-C2'-C1'	-7.22	95.72	101.50
1	AA	438	G	P-O3'-C3'	7.21	128.35	119.70
1	CA	438	G	P-O3'-C3'	7.20	128.33	119.70
23	DA	1022	G	P-O3'-C3'	7.13	128.26	119.70
23	BA	242	G	C3'-C2'-C1'	-7.05	95.86	101.50
1	AA	1529	G	C1'-O4'-C4'	-7.05	104.26	109.90
23	DA	242	G	C3'-C2'-C1'	-7.02	95.88	101.50
1	AA	136(B)	C	O3'-P-O5'	-6.97	90.75	104.00
23	BA	1496	A	O4'-C1'-N9	6.94	113.75	108.20
1	CA	1129	C	P-O3'-C3'	6.94	128.03	119.70
23	DA	1395	A	C1'-O4'-C4'	-6.94	104.35	109.90
23	BA	332	A	P-O3'-C3'	6.92	128.00	119.70
1	AA	1129	C	P-O3'-C3'	6.89	127.97	119.70
23	DA	1698	A	C3'-C2'-C1'	-6.84	96.03	101.50
23	DA	989	G	C1'-O4'-C4'	-6.84	104.43	109.90
23	BA	1395	A	C1'-O4'-C4'	-6.82	104.44	109.90
23	DA	332	A	P-O3'-C3'	6.82	127.88	119.70
23	BA	1558	A	P-O3'-C3'	6.80	127.86	119.70
1	AA	1201	A	P-O3'-C3'	6.80	127.86	119.70
1	AA	119	A	P-O3'-C3'	6.78	127.84	119.70
23	BA	2506	U	O4'-C1'-N1	6.75	113.60	108.20
1	AA	687	A	P-O3'-C3'	6.74	127.79	119.70
23	BA	1542	G	P-O3'-C3'	6.70	127.75	119.70
23	BA	989	G	C1'-O4'-C4'	-6.69	104.55	109.90
23	BA	1984	G	C4'-C3'-C2'	-6.67	95.93	102.60
1	CA	687	A	P-O3'-C3'	6.64	127.66	119.70
23	DA	1545	A	C1'-O4'-C4'	-6.63	104.60	109.90
23	DA	131	G	C1'-O4'-C4'	-6.62	104.60	109.90
23	BA	1698	A	C3'-C2'-C1'	-6.59	96.22	101.50
23	BA	2445	G	C4'-C3'-C2'	-6.58	96.02	102.60
23	BA	1022	G	P-O3'-C3'	6.58	127.59	119.70
23	DA	1558	A	P-O3'-C3'	6.57	127.59	119.70
1	AA	366	C	P-O3'-C3'	6.56	127.58	119.70
23	BA	1021	A	C3'-C2'-C1'	-6.56	96.25	101.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
23	DA	2506	U	O4'-C1'-N1	6.54	113.44	108.20
23	DA	1542	G	P-O3'-C3'	6.54	127.55	119.70
1	AA	1300	G	P-O3'-C3'	6.52	127.52	119.70
23	DA	1266	G	C3'-C2'-C1'	-6.50	96.30	101.50
1	CA	1300	G	P-O3'-C3'	6.50	127.50	119.70
23	BA	1046	A	C5-C6-N1	-6.48	114.46	117.70
23	DA	372	G	O4'-C1'-N9	6.48	113.38	108.20
23	DA	2049	G	C1'-O4'-C4'	-6.47	104.72	109.90
1	CA	1201	A	P-O3'-C3'	6.47	127.47	119.70
23	DA	2445	G	C4'-C3'-C2'	-6.47	96.13	102.60
23	BA	1379	A	P-O3'-C3'	6.47	127.46	119.70
23	DA	2447	G	P-O3'-C3'	6.47	127.46	119.70
23	DA	783	A	C3'-C2'-C1'	-6.45	96.34	101.50
1	CA	366	C	P-O3'-C3'	6.42	127.41	119.70
1	CA	358	U	P-O3'-C3'	-6.39	112.03	119.70
23	DA	1427	A	P-O3'-C3'	6.39	127.37	119.70
23	DA	933	A	C4'-C3'-C2'	-6.38	96.22	102.60
23	DA	1385	G	C3'-C2'-C1'	-6.37	96.41	101.50
1	CA	119	A	P-O3'-C3'	6.36	127.33	119.70
23	DA	1698	A	C1'-O4'-C4'	-6.36	104.81	109.90
23	DA	783	A	C2-N3-C4	-6.35	107.43	110.60
23	BA	2049	G	C1'-O4'-C4'	-6.33	104.84	109.90
1	AA	824	C	C3'-C2'-C1'	-6.32	96.44	101.50
23	BA	933	A	C4'-C3'-C2'	-6.32	96.28	102.60
23	DA	1608	A	C1'-O4'-C4'	-6.30	104.86	109.90
23	BA	1545	A	C1'-O4'-C4'	-6.29	104.87	109.90
23	DA	761	A	C6-C5-N7	-6.29	127.90	132.30
23	BA	210	C	C6-N1-C2	6.27	122.81	120.30
23	BA	1786	A	N9-C1'-C2'	6.24	122.11	114.00
23	BA	131	G	C1'-O4'-C4'	-6.24	104.91	109.90
23	DA	933	A	C1'-O4'-C4'	-6.24	104.91	109.90
23	DA	1984	G	C4'-C3'-C2'	-6.23	96.37	102.60
23	BA	2225	A	P-O3'-C3'	6.22	127.17	119.70
1	CA	1067	A	P-O3'-C3'	6.22	127.16	119.70
23	BA	791	C	C3'-C2'-C1'	-6.21	96.53	101.50
23	BA	1385	G	C3'-C2'-C1'	-6.20	96.54	101.50
23	DA	131	G	O4'-C4'-C3'	-6.19	97.81	104.00
23	DA	1046	A	C5-C6-N1	-6.19	114.60	117.70
23	BA	859	G	C3'-C2'-C1'	-6.19	96.55	101.50
23	BA	2613	U	C3'-C2'-C1'	-6.19	96.55	101.50
23	DA	1496	A	O4'-C1'-N9	6.17	113.14	108.20
1	CA	246	A	C3'-C2'-C1'	-6.17	96.56	101.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
23	DA	2225	A	P-O3'-C3'	6.16	127.09	119.70
1	CA	1529	G	O4'-C1'-N9	6.15	113.12	108.20
1	AA	1067	A	P-O3'-C3'	6.14	127.07	119.70
23	BA	1761	C	C1'-O4'-C4'	-6.14	104.99	109.90
23	DA	859	G	C3'-C2'-C1'	-6.14	96.59	101.50
23	DA	1379	A	P-O3'-C3'	6.13	127.05	119.70
1	AA	136(B)	C	P-O3'-C3'	-6.12	112.35	119.70
23	BA	2447	G	P-O3'-C3'	6.12	127.05	119.70
23	DA	450	G	C5-C6-N1	-6.09	108.45	111.50
23	BA	1266	G	C3'-C2'-C1'	-6.09	96.62	101.50
23	DA	791	C	C3'-C2'-C1'	-6.08	96.63	101.50
23	DA	1021	A	C3'-C2'-C1'	-6.08	96.63	101.50
23	DA	673	C	C4'-C3'-C2'	-6.08	96.52	102.60
1	CA	824	C	C3'-C2'-C1'	-6.07	96.64	101.50
1	AA	1529	G	O4'-C1'-N9	6.04	113.03	108.20
23	DA	1786	A	N9-C1'-C2'	6.04	121.85	114.00
23	DA	2033	A	P-O3'-C3'	6.04	126.94	119.70
23	BA	372	G	O4'-C1'-N9	6.03	113.03	108.20
23	DA	783	A	N1-C6-N6	6.02	122.21	118.60
23	DA	760	G	C4'-C3'-C2'	-6.01	96.59	102.60
23	DA	2613	U	C3'-C2'-C1'	-6.00	96.70	101.50
23	BA	783	A	N1-C6-N6	5.98	122.19	118.60
23	BA	989	G	O4'-C1'-N9	5.98	112.98	108.20
23	BA	2052	G	C4'-C3'-C2'	-5.97	96.63	102.60
23	BA	2056	G	N9-C4-C5	5.96	107.78	105.40
23	DA	686	G	O4'-C1'-N9	5.96	112.96	108.20
23	BA	933	A	C1'-O4'-C4'	-5.93	105.15	109.90
23	DA	2481	G	P-O3'-C3'	5.93	126.82	119.70
1	AA	1504	G	P-O3'-C3'	5.91	126.79	119.70
1	CA	1285	A	P-O3'-C3'	5.90	126.78	119.70
23	BA	938	G	C4'-C3'-C2'	-5.90	96.70	102.60
23	DA	2562	U	C4'-C3'-C2'	-5.89	96.70	102.60
1	AA	1285	A	P-O3'-C3'	5.89	126.77	119.70
23	BA	2033	A	P-O3'-C3'	5.89	126.76	119.70
23	BA	825	C	C4'-C3'-C2'	-5.87	96.73	102.60
23	BA	407	G	C4'-C3'-C2'	-5.86	96.74	102.60
1	AA	246	A	C3'-C2'-C1'	-5.85	96.82	101.50
22	CV	6188	C	P-O3'-C3'	5.83	126.70	119.70
23	BA	783	A	C3'-C2'-C1'	-5.82	96.84	101.50
23	BA	1427	A	P-O3'-C3'	5.82	126.68	119.70
22	AV	6188	C	P-O3'-C3'	5.82	126.68	119.70
23	BA	1786	A	O4'-C1'-C2'	-5.82	99.98	105.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	810	C	C1'-O4'-C4'	-5.79	105.26	109.90
23	DA	2049	G	C4'-C3'-C2'	-5.79	96.81	102.60
23	BA	1698	A	C1'-O4'-C4'	-5.78	105.28	109.90
23	DA	825	C	C4'-C3'-C2'	-5.78	96.83	102.60
23	BA	444	C	C1'-O4'-C4'	-5.77	105.28	109.90
23	BA	2428	G	P-O3'-C3'	5.76	126.61	119.70
23	DA	807	U	C4'-C3'-C2'	-5.76	96.84	102.60
23	BA	807	U	C4'-C3'-C2'	-5.76	96.84	102.60
1	AA	913	A	P-O3'-C3'	5.75	126.59	119.70
23	DA	2098	U	P-O3'-C3'	5.72	126.56	119.70
23	BA	2278	A	C1'-O4'-C4'	-5.70	105.34	109.90
1	CA	913	A	P-O3'-C3'	5.70	126.54	119.70
23	DA	733	G	O4'-C1'-N9	-5.70	103.64	108.20
1	AA	7	G	C1'-O4'-C4'	-5.69	105.35	109.90
1	CA	7	G	C1'-O4'-C4'	-5.69	105.34	109.90
23	BA	2883	A	O4'-C1'-N9	5.69	112.75	108.20
22	CV	6171	U	P-O3'-C3'	5.69	126.53	119.70
23	DA	1761	C	C1'-O4'-C4'	-5.67	105.36	109.90
23	DA	444	C	C4'-C3'-C2'	-5.66	96.94	102.60
23	DA	761	A	N7-C8-N9	5.66	116.63	113.80
22	AV	6171	U	P-O3'-C3'	5.65	126.48	119.70
23	DA	680	G	O4'-C1'-N9	-5.65	103.68	108.20
23	DA	2512	C	C4'-C3'-C2'	-5.65	96.95	102.60
23	BA	974(B)	C	C3'-C2'-C1'	-5.64	96.99	101.50
23	BA	2098	U	P-O3'-C3'	5.63	126.46	119.70
23	DA	2056	G	N3-C4-N9	-5.63	122.62	126.00
23	BA	761	A	N7-C8-N9	5.61	116.61	113.80
23	BA	933	A	O4'-C4'-C3'	-5.60	98.40	104.00
23	BA	1608	A	C1'-O4'-C4'	-5.60	105.42	109.90
23	BA	2481	G	P-O3'-C3'	5.59	126.41	119.70
23	DA	2278	A	C1'-O4'-C4'	-5.58	105.44	109.90
23	DA	783	A	C5-N7-C8	-5.58	101.11	103.90
23	BA	131	G	O4'-C4'-C3'	-5.57	98.43	104.00
24	DB	66	A	P-O3'-C3'	5.57	126.39	119.70
23	DA	2428	G	P-O3'-C3'	5.57	126.39	119.70
1	AA	890	G	O4'-C1'-N9	5.56	112.65	108.20
23	BA	2689	U	P-O3'-C3'	5.56	126.37	119.70
23	DA	989	G	O4'-C1'-N9	5.53	112.62	108.20
23	BA	940	G	C4'-C3'-C2'	-5.51	97.09	102.60
23	DA	974(B)	C	C3'-C2'-C1'	-5.51	97.10	101.50
23	BA	226	G	C3'-C2'-C1'	-5.50	97.10	101.50
23	DA	95	G	C4'-C3'-C2'	-5.50	97.11	102.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
24	BB	66	A	P-O3'-C3'	5.49	126.29	119.70
23	DA	1255	U	C1'-O4'-C4'	-5.49	105.51	109.90
23	DA	940	G	C4'-C3'-C2'	-5.48	97.12	102.60
23	BA	2562	U	C4'-C3'-C2'	-5.48	97.12	102.60
23	DA	450	G	C4-C5-C6	5.47	122.08	118.80
23	DA	85	G	C3'-C2'-C1'	-5.46	97.13	101.50
1	AA	356	A	C4'-C3'-C2'	-5.45	97.15	102.60
23	DA	1275	A	C3'-C2'-C1'	-5.45	97.14	101.50
1	CA	1504	G	P-O3'-C3'	5.44	126.23	119.70
23	DA	938	G	C4'-C3'-C2'	-5.43	97.17	102.60
23	BA	760	G	C4'-C3'-C2'	-5.42	97.18	102.60
1	AA	358	U	P-O3'-C3'	-5.41	113.21	119.70
23	BA	95	G	C4'-C3'-C2'	-5.40	97.20	102.60
23	DA	527	C	O4'-C1'-N1	5.40	112.52	108.20
23	DA	407	G	C4'-C3'-C2'	-5.39	97.21	102.60
23	BA	761	A	C6-C5-N7	-5.39	128.53	132.30
23	DA	479	A	C1'-O4'-C4'	-5.39	105.59	109.90
23	BA	673	C	C4'-C3'-C2'	-5.39	97.21	102.60
23	DA	1786	A	O4'-C1'-C2'	-5.38	100.42	105.80
23	BA	1255	U	C1'-O4'-C4'	-5.38	105.60	109.90
23	DA	2052	G	C4'-C3'-C2'	-5.37	97.23	102.60
23	BA	1162	G	C4'-C3'-C2'	-5.37	97.23	102.60
1	CA	890	G	O4'-C1'-N9	5.37	112.49	108.20
23	BA	527	C	O4'-C1'-N1	5.36	112.49	108.20
23	BA	2681	C	C3'-C2'-C1'	-5.36	97.21	101.50
23	BA	1022	G	C3'-C2'-C1'	5.36	105.79	101.50
23	DA	226	G	C3'-C2'-C1'	-5.35	97.22	101.50
23	DA	2789	C	C1'-O4'-C4'	-5.35	105.62	109.90
1	AA	328	C	P-O3'-C3'	5.35	126.12	119.70
23	BA	2549	G	C4'-C3'-C2'	-5.35	97.25	102.60
23	BA	2512	C	C4'-C3'-C2'	-5.34	97.26	102.60
1	AA	890	G	C1'-O4'-C4'	-5.34	105.63	109.90
23	BA	956	G	C1'-O4'-C4'	-5.32	105.64	109.90
1	AA	1064	G	P-O3'-C3'	5.32	126.08	119.70
1	AA	890	G	C3'-C2'-C1'	-5.31	97.25	101.50
23	BA	84	A	C3'-C2'-C1'	-5.30	97.26	101.50
23	DA	761	A	C5-C6-N1	-5.30	115.05	117.70
23	BA	372	G	C3'-C2'-C1'	-5.29	97.27	101.50
23	DA	2689	U	P-O3'-C3'	5.29	126.05	119.70
23	DA	1162	G	C4'-C3'-C2'	-5.29	97.31	102.60
23	DA	1616	A	O4'-C1'-N9	5.28	112.43	108.20
23	BA	2049	G	C4'-C3'-C2'	-5.28	97.32	102.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
23	BA	2056	G	C4'-C3'-C2'	-5.28	97.32	102.60
23	BA	2818	G	O4'-C4'-C3'	-5.28	98.72	104.00
23	DA	210	C	C6-N1-C2	5.27	122.41	120.30
23	BA	1936	A	P-O3'-C3'	5.27	126.02	119.70
23	BA	807	U	O4'-C4'-C3'	-5.27	98.73	104.00
23	BA	444	C	C4'-C3'-C2'	-5.26	97.34	102.60
23	BA	479	A	C1'-O4'-C4'	-5.25	105.70	109.90
1	CA	1525	G	C4'-C3'-C2'	-5.25	97.35	102.60
1	AA	1525	G	C4'-C3'-C2'	-5.24	97.36	102.60
23	DA	532	A	C1'-O4'-C4'	-5.24	105.71	109.90
1	AA	428	G	P-O3'-C3'	5.22	125.97	119.70
23	BA	214	G	O4'-C1'-N9	5.22	112.38	108.20
23	DA	1613	G	C4'-C3'-C2'	-5.21	97.39	102.60
1	CA	890	G	C3'-C2'-C1'	-5.21	97.33	101.50
23	BA	1613	G	C4'-C3'-C2'	-5.21	97.39	102.60
23	DA	974(B)	C	O4'-C1'-C2'	-5.21	100.59	105.80
23	DA	1855	G	C4'-C3'-C2'	-5.20	97.40	102.60
23	BA	601	C	C4'-C3'-C2'	-5.18	97.42	102.60
23	DA	1012	U	C3'-C2'-C1'	-5.18	97.36	101.50
23	DA	444	C	C1'-O4'-C4'	-5.17	105.77	109.90
23	BA	532	A	C1'-O4'-C4'	-5.15	105.78	109.90
23	BA	2447	G	C3'-C2'-C1'	-5.14	97.38	101.50
1	CA	527	G	C4'-C3'-C2'	-5.14	97.46	102.60
23	DA	784	A	C3'-C2'-C1'	5.14	105.61	101.50
1	AA	1077	G	C4'-C3'-C2'	-5.13	97.47	102.60
1	CA	810	C	C1'-O4'-C4'	-5.13	105.79	109.90
23	DA	1559	G	C1'-O4'-C4'	-5.13	105.79	109.90
23	DA	2056	G	C4'-C3'-C2'	-5.13	97.47	102.60
23	BA	2789	C	C1'-O4'-C4'	-5.13	105.80	109.90
23	DA	208	C	C6-N1-C2	5.13	122.35	120.30
23	DA	692	C	C4'-C3'-C2'	-5.13	97.47	102.60
23	DA	463	G	C4'-C3'-C2'	-5.12	97.48	102.60
23	DA	807	U	O4'-C4'-C3'	-5.12	98.88	104.00
23	BA	1929	G	C3'-C2'-C1'	-5.12	97.40	101.50
1	CA	356	A	C4'-C3'-C2'	-5.12	97.48	102.60
1	CA	890	G	C1'-O4'-C4'	-5.12	105.80	109.90
23	DA	2744	G	C4'-C3'-C2'	-5.12	97.48	102.60
23	DA	1504	C	N1-C1'-C2'	-5.11	106.37	112.00
23	DA	933	A	O4'-C4'-C3'	-5.11	98.89	104.00
23	DA	1267	U	P-O3'-C3'	5.11	125.83	119.70
1	AA	1361	G	OP2-P-O3'	5.11	116.43	105.20
23	BA	974(B)	C	C1'-O4'-C4'	-5.09	105.82	109.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
23	DA	2681	C	C3'-C2'-C1'	-5.09	97.43	101.50
23	DA	1496	A	C1'-O4'-C4'	-5.09	105.83	109.90
23	BA	692	C	C4'-C3'-C2'	-5.08	97.52	102.60
23	BA	2744	G	C4'-C3'-C2'	-5.07	97.53	102.60
1	CA	1077	G	C4'-C3'-C2'	-5.07	97.53	102.60
1	CA	1064	G	P-O3'-C3'	5.07	125.78	119.70
23	DA	738	G	C1'-O4'-C4'	-5.06	105.85	109.90
1	CA	250	A	P-O3'-C3'	5.06	125.77	119.70
1	CA	792	A	C3'-C2'-C1'	-5.06	97.45	101.50
24	DB	42	C	C1'-O4'-C4'	-5.06	105.85	109.90
23	DA	1022	G	C3'-C2'-C1'	5.05	105.54	101.50
23	DA	324	A	C4'-C3'-C2'	-5.04	97.56	102.60
23	BA	265	A	C1'-O4'-C4'	-5.04	105.87	109.90
23	BA	1494	A	P-O3'-C3'	5.04	125.74	119.70
23	BA	1189	A	O3'-P-O5'	-5.03	94.44	104.00
23	BA	741	G	C4'-C3'-C2'	-5.03	97.57	102.60
23	DA	1929	G	C3'-C2'-C1'	-5.03	97.48	101.50
23	DA	672	C	C4'-C3'-C2'	-5.02	97.58	102.60
1	CA	328	C	P-O3'-C3'	5.02	125.72	119.70
23	BA	671	C	C4'-C3'-C2'	-5.02	97.58	102.60
23	BA	85	G	C3'-C2'-C1'	-5.02	97.49	101.50
1	CA	1027	C	P-O3'-C3'	-5.01	113.68	119.70
23	BA	35	G	C4'-C3'-C2'	-5.01	97.59	102.60
23	BA	805	G	O4'-C1'-N9	5.01	112.21	108.20
23	DA	1743	G	C4'-C3'-C2'	-5.01	97.59	102.60
23	BA	1504	C	N1-C1'-C2'	-5.00	106.50	112.00
23	BA	1654	A	C4'-C3'-C2'	-5.00	97.60	102.60

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
28	BF	75	LYS	Peptide
34	BL	52	GLU	Peptide
28	DF	75	LYS	Peptide
34	DL	52	GLU	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	AA	32372	0	16339	682	0
1	CA	32372	0	16339	694	0
2	AB	1901	0	1951	103	0
2	CB	1901	0	1951	103	0
3	AC	1613	0	1677	95	0
3	CC	1613	0	1677	95	0
4	AD	1703	0	1764	100	0
4	CD	1703	0	1765	92	0
5	AE	1156	0	1213	68	0
5	CE	1156	0	1213	66	0
6	AF	843	0	857	46	0
6	CF	843	0	857	46	0
7	AG	1257	0	1296	60	0
7	CG	1257	0	1296	62	0
8	AH	1116	0	1177	57	0
8	CH	1116	0	1177	59	0
9	AI	1011	0	1043	62	0
9	CI	1011	0	1043	58	0
10	AJ	795	0	840	59	0
10	CJ	795	0	840	59	0
11	AK	885	0	904	56	0
11	CK	885	0	904	50	0
12	AL	971	0	1057	60	0
12	CL	971	0	1057	66	0
13	AM	929	0	987	66	0
13	CM	929	0	987	64	0
14	AN	492	0	529	29	0
14	CN	492	0	533	26	0
15	AO	734	0	771	26	0
15	CO	734	0	771	29	0
16	AP	701	0	720	30	0
16	CP	701	0	720	38	0
17	AQ	824	0	893	42	0
17	CQ	824	0	893	44	0
18	AR	574	0	644	30	0
18	CR	574	0	644	32	0
19	AS	630	0	652	59	0
19	CS	630	0	652	54	0
20	AT	762	0	859	32	0
20	CT	762	0	859	31	0
21	AU	209	0	221	13	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
21	CU	209	0	221	14	0
22	AV	736	0	378	19	0
22	CV	736	0	378	18	0
23	BA	59442	0	29966	1292	0
23	DA	59442	0	29966	1295	0
24	BB	2551	0	1295	53	0
24	DB	2551	0	1295	53	0
25	BC	2105	0	2182	202	0
25	DC	2105	0	2182	214	0
26	BD	1564	0	1629	117	0
26	DD	1564	0	1629	112	0
27	BE	1587	0	1632	107	0
27	DE	1587	0	1632	106	0
28	BF	1475	0	1537	109	0
28	DF	1475	0	1537	109	0
29	BG	1223	0	1282	64	0
29	DG	1223	0	1282	66	0
30	BH	1133	0	1220	83	0
30	DH	1133	0	1220	80	0
31	BI	254	0	275	8	0
31	DI	254	0	275	8	0
32	BJ	1097	0	1168	82	0
32	DJ	1097	0	1168	83	0
33	BK	932	0	994	54	0
33	DK	932	0	994	56	0
34	BL	1114	0	1187	167	0
34	DL	1114	0	1187	169	0
35	BM	1079	0	1127	96	0
35	DM	1079	0	1127	99	0
36	BN	960	0	1021	73	0
36	DN	960	0	1021	71	0
37	BO	771	0	832	67	0
37	DO	771	0	832	70	0
38	BP	1144	0	1211	69	0
38	DP	1144	0	1211	74	0
39	BQ	953	0	1013	71	0
39	DQ	953	0	1013	71	0
40	BR	779	0	852	76	0
40	DR	779	0	852	73	0
41	BS	891	0	951	50	0
41	DS	891	0	951	52	0
42	BT	726	0	778	58	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
42	DT	726	0	778	59	0
43	BU	776	0	870	78	0
43	DU	776	0	870	76	0
44	BV	1492	0	1513	94	0
44	DV	1492	0	1513	92	0
45	BW	605	0	628	36	0
45	DW	605	0	628	38	0
46	BX	695	0	764	69	0
46	DX	695	0	764	68	0
47	BY	521	0	575	45	0
47	DY	521	0	575	45	0
48	BZ	468	0	523	20	0
48	DZ	468	0	523	24	0
49	B1	226	0	225	17	0
49	D1	226	0	225	17	0
50	B2	405	0	420	27	0
50	D2	405	0	420	29	0
51	B3	381	0	391	28	0
51	D3	381	0	391	26	0
52	B4	419	0	467	30	0
52	D4	419	0	467	32	0
53	B5	508	0	576	54	0
53	D5	508	0	576	55	0
54	AA	279	0	0	0	0
54	AB	3	0	0	0	0
54	AC	2	0	0	0	0
54	AD	3	0	0	0	0
54	AE	5	0	0	0	0
54	AF	1	0	0	0	0
54	AG	2	0	0	0	0
54	AH	1	0	0	0	0
54	AI	1	0	0	0	0
54	AK	2	0	0	0	0
54	AL	2	0	0	0	0
54	AM	1	0	0	0	0
54	AN	1	0	0	0	0
54	AO	3	0	0	0	0
54	AQ	3	0	0	0	0
54	AS	1	0	0	0	0
54	AT	2	0	0	0	0
54	AV	8	0	0	0	0
54	B2	3	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
54	B3	1	0	0	0	0
54	B4	1	0	0	0	0
54	B5	1	0	0	0	0
54	BA	781	0	0	0	0
54	BB	27	0	0	0	0
54	BC	3	0	0	0	0
54	BE	2	0	0	0	0
54	BF	1	0	0	0	0
54	BH	1	0	0	0	0
54	BI	1	0	0	0	0
54	BJ	3	0	0	0	0
54	BK	1	0	0	0	0
54	BM	3	0	0	0	0
54	BN	2	0	0	0	0
54	BO	1	0	0	0	0
54	BP	2	0	0	0	0
54	BR	1	0	0	0	0
54	BS	1	0	0	0	0
54	BT	2	0	0	0	0
54	BU	2	0	0	0	0
54	BV	3	0	0	0	0
54	BW	1	0	0	0	0
54	BX	2	0	0	0	0
54	BY	3	0	0	0	0
54	CA	372	0	0	0	0
54	CB	3	0	0	0	0
54	CC	8	0	0	0	0
54	CD	3	0	0	0	0
54	CE	5	0	0	0	0
54	CF	2	0	0	0	0
54	CG	3	0	0	0	0
54	CK	1	0	0	0	0
54	CL	6	0	0	0	0
54	CM	2	0	0	0	0
54	CN	1	0	0	0	0
54	CO	4	0	0	0	0
54	CQ	2	0	0	0	0
54	CV	7	0	0	0	0
54	D1	1	0	0	0	0
54	D2	3	0	0	0	0
54	D3	1	0	0	0	0
54	D4	4	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
54	D5	3	0	0	0	0
54	DA	964	0	0	0	0
54	DB	43	0	0	0	0
54	DC	3	0	0	0	0
54	DD	6	0	0	0	0
54	DE	3	0	0	0	0
54	DF	2	0	0	0	0
54	DG	3	0	0	0	0
54	DI	1	0	0	0	0
54	DJ	6	0	0	0	0
54	DK	3	0	0	0	0
54	DL	3	0	0	0	0
54	DM	1	0	0	0	0
54	DN	1	0	0	0	0
54	DO	1	0	0	0	0
54	DP	2	0	0	0	0
54	DQ	2	0	0	0	0
54	DS	2	0	0	0	0
54	DU	1	0	0	0	0
54	DV	4	0	0	0	0
54	DW	1	0	0	0	0
54	DX	3	0	0	0	0
54	DY	2	0	0	0	0
55	AD	1	0	0	0	0
55	AN	1	0	0	0	0
55	CD	1	0	0	0	0
55	CN	1	0	0	0	0
All	All	283641	0	191757	9280	0

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 20.

All (9280) 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)
1:AA:1399:C:H4'	1:AA:1400:C:H5'	1.26	1.16
34:BL:33:ARG:HG3	34:BL:36:LYS:HD3	1.31	1.12
34:BL:128:HIS:HA	34:BL:147:LEU:HB3	1.30	1.11
23:DA:2015:A:H1'	50:D2:2:ALA:HA	1.33	1.08
34:DL:33:ARG:HG3	34:DL:36:LYS:HD3	1.33	1.08
1:CA:1399:C:H4'	1:CA:1400:C:H5'	1.27	1.08

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BA:2015:A:H1'	50:B2:2:ALA:HA	1.34	1.07
34:DL:49:ARG:HG2	34:DL:50:ARG:H	1.18	1.06
34:DL:128:HIS:HA	34:DL:147:LEU:HB3	1.30	1.06
34:BL:49:ARG:HG2	34:BL:50:ARG:H	1.16	1.04
7:AG:113:GLU:HB2	7:AG:119:ARG:HG2	1.38	1.04
45:BW:23:VAL:HA	45:BW:38:VAL:HG22	1.37	1.03
32:DJ:70:ALA:HB2	32:DJ:135:LEU:HD12	1.41	1.03
45:DW:23:VAL:HA	45:DW:38:VAL:HG22	1.37	1.03
37:BO:24:LEU:HD12	37:BO:84:GLN:HB3	1.38	1.03
30:BH:92:VAL:HG13	30:BH:120:ILE:HB	1.41	1.03
7:CG:113:GLU:HB2	7:CG:119:ARG:HG2	1.38	1.02
9:AI:19:LEU:HD21	9:AI:59:PHE:HB3	1.41	1.02
37:DO:24:LEU:HD12	37:DO:84:GLN:HB3	1.38	1.02
39:DQ:92:ARG:HH11	39:DQ:92:ARG:HB2	1.24	1.01
39:BQ:92:ARG:HB2	39:BQ:92:ARG:HH11	1.25	1.01
25:BC:242:ARG:H	25:BC:242:ARG:HD3	1.21	1.00
9:CI:19:LEU:HD21	9:CI:59:PHE:HB3	1.40	1.00
15:AO:63:ARG:HH21	15:AO:87:ILE:HG21	1.27	0.99
25:DC:242:ARG:HD3	25:DC:242:ARG:H	1.23	0.99
5:AE:76:ILE:HG12	5:AE:77:PRO:HD2	1.43	0.99
30:DH:92:VAL:HG13	30:DH:120:ILE:HB	1.42	0.99
43:DU:8:LYS:H	43:DU:8:LYS:HZ2	1.07	0.98
23:DA:2681:C:H5	23:DA:2725:A:H62	0.99	0.98
23:DA:2781:A:H5'	23:DA:2782:G:H5'	1.46	0.98
32:BJ:70:ALA:HB2	32:BJ:135:LEU:HD12	1.41	0.98
42:BT:11:PRO:HA	42:BT:28:PHE:HB3	1.47	0.97
23:BA:2781:A:H5'	23:BA:2782:G:H5'	1.44	0.97
23:DA:1163:G:H2'	23:DA:1164:G:H5''	1.46	0.97
15:CO:63:ARG:HH21	15:CO:87:ILE:HG21	1.28	0.96
23:BA:1163:G:H2'	23:BA:1164:G:H5''	1.46	0.96
34:BL:50:ARG:HB2	53:B5:60:LEU:HD11	1.47	0.96
13:CM:76:ALA:HA	13:CM:79:LYS:HE2	1.48	0.96
5:CE:76:ILE:HG12	5:CE:77:PRO:HD2	1.43	0.95
42:DT:11:PRO:HA	42:DT:28:PHE:HB3	1.47	0.95
1:AA:1443:G:H3'	1:AA:1446:A:H5''	1.48	0.95
23:BA:1021:A:H8	23:BA:1021:A:H3'	1.31	0.95
32:BJ:157:ARG:H	32:BJ:158:PRO:HD3	1.30	0.95
23:DA:2068:U:H3	23:DA:2430:A:H2	1.13	0.95
1:CA:522:C:H41	12:CL:52:ARG:HH22	1.11	0.95
2:CB:185:ILE:HG22	2:CB:199:TYR:HB2	1.47	0.95
1:CA:1443:G:H3'	1:CA:1446:A:H5''	1.48	0.95

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
44:DV:18:LEU:HD23	44:DV:25:PRO:HG3	1.49	0.95
1:AA:955:U:H1'	1:AA:1227:A:H61	1.31	0.94
23:DA:780:G:H21	23:DA:783:A:H62	1.15	0.94
23:DA:1899:G:H21	23:DA:1902:C:H41	1.07	0.94
2:AB:185:ILE:HG22	2:AB:199:TYR:HB2	1.47	0.94
23:BA:2420:C:OP1	53:B5:34:TRP:HA	1.68	0.93
32:DJ:42:GLU:HA	32:DJ:82:LYS:HB3	1.50	0.93
23:BA:1021:A:H3'	23:BA:1021:A:C8	2.04	0.93
23:BA:1899:G:H21	23:BA:1902:C:H41	1.08	0.93
10:CJ:50:ILE:HB	14:CN:41:ARG:HH21	1.33	0.93
1:CA:955:U:H1'	1:CA:1227:A:H61	1.31	0.93
1:CA:1378:C:H5''	7:CG:6:ARG:HE	1.33	0.92
29:BG:16:SER:HB2	29:BG:27:LYS:HB2	1.51	0.92
32:BJ:42:GLU:HA	32:BJ:82:LYS:HB3	1.50	0.92
23:DA:1046:A:N3	31:DI:4:LYS:HD3	1.84	0.92
13:AM:76:ALA:HA	13:AM:79:LYS:HE2	1.48	0.92
35:DM:22:LYS:HA	35:DM:22:LYS:HE2	1.50	0.92
32:DJ:157:ARG:H	32:DJ:158:PRO:HD3	1.30	0.92
35:BM:22:LYS:HA	35:BM:22:LYS:HE2	1.51	0.92
28:BF:60:LEU:HD11	28:BF:92:VAL:HG11	1.52	0.92
23:BA:1019:U:HO2'	23:BA:1021:A:H2	0.99	0.91
23:DA:1021:A:C8	23:DA:1021:A:H3'	2.05	0.91
34:DL:50:ARG:HB2	53:D5:60:LEU:HD11	1.49	0.91
28:DF:60:LEU:HD11	28:DF:92:VAL:HG11	1.52	0.91
34:DL:49:ARG:HG2	34:DL:50:ARG:N	1.84	0.91
23:DA:1019:U:HO2'	23:DA:1021:A:H2	0.95	0.91
23:BA:886:C:H2'	23:BA:887:A:H4'	1.52	0.91
23:BA:2068:U:H3	23:BA:2430:A:H2	1.11	0.91
34:DL:114:ILE:HD11	34:DL:130:PHE:CD1	2.05	0.91
23:DA:2420:C:OP1	53:D5:34:TRP:HA	1.71	0.90
23:DA:587:C:H42	34:DL:33:ARG:HG2	1.35	0.90
29:DG:16:SER:HB2	29:DG:27:LYS:HB2	1.52	0.90
12:CL:26:LEU:HD13	12:CL:27:LYS:H	1.36	0.90
10:AJ:50:ILE:HB	14:AN:41:ARG:HH21	1.34	0.90
23:BA:1046:A:N3	31:BI:4:LYS:HD3	1.84	0.90
23:BA:587:C:H42	34:BL:33:ARG:HG2	1.34	0.90
23:BA:2681:C:H5	23:BA:2725:A:H62	0.98	0.90
33:DK:3:GLN:HB2	33:DK:4:PRO:HD2	1.53	0.90
44:BV:18:LEU:HD23	44:BV:25:PRO:HG3	1.52	0.90
23:BA:761:A:H8	23:BA:761:A:O5'	1.53	0.90
47:DY:2:LYS:HA	47:DY:5:GLU:CD	1.90	0.90

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DA:547:A:H2'	23:DA:548:A:C8	2.07	0.90
23:DA:1021:A:H3'	23:DA:1021:A:H8	1.32	0.90
19:AS:19:VAL:HG21	19:AS:44:MET:HG3	1.55	0.89
23:DA:2426:A:H3'	23:DA:2427:C:H5''	1.54	0.89
43:DU:81:LYS:HD2	43:DU:96:ILE:HD12	1.54	0.89
1:CA:1223:C:H5'	1:CA:1224:G:H5''	1.54	0.89
1:CA:1261:A:H62	1:CA:1274:G:H21	1.18	0.89
21:CU:22:ARG:HD2	21:CU:23:PRO:HD2	1.55	0.89
23:DA:761:A:O5'	23:DA:761:A:H8	1.56	0.89
23:BA:996:A:H4'	39:BQ:92:ARG:NH1	1.87	0.89
41:BS:24:ILE:HG21	41:BS:36:LEU:HD21	1.55	0.89
1:CA:736:C:H2'	1:CA:737:A:C8	2.07	0.89
23:BA:547:A:H2'	23:BA:548:A:C8	2.08	0.89
12:AL:26:LEU:HD13	12:AL:27:LYS:H	1.36	0.89
49:D1:50:THR:HG22	49:D1:51:TYR:H	1.37	0.89
1:AA:134:A:H61	16:AP:25:ARG:NH1	1.69	0.89
47:BY:2:LYS:HA	47:BY:5:GLU:CD	1.94	0.89
1:AA:736:C:H2'	1:AA:737:A:C8	2.08	0.89
19:CS:19:VAL:HG21	19:CS:44:MET:HG3	1.54	0.88
23:DA:996:A:H4'	39:DQ:92:ARG:NH1	1.87	0.88
23:BA:142:G:H4'	42:BT:35:THR:HG21	1.55	0.88
29:DG:101:ARG:HE	29:DG:101:ARG:H	1.18	0.88
1:CA:979:C:H3'	1:CA:980:C:H5''	1.55	0.88
34:BL:49:ARG:HG2	34:BL:50:ARG:N	1.83	0.88
34:BL:114:ILE:HD11	34:BL:130:PHE:CD1	2.07	0.88
23:BA:780:G:H21	23:BA:783:A:H62	1.14	0.88
23:DA:954:G:H5''	35:DM:13:GLN:HG3	1.54	0.88
23:DA:886:C:H2'	23:DA:887:A:H4'	1.53	0.88
5:AE:51:VAL:HB	5:AE:52:PRO:HD3	1.56	0.88
23:DA:2331:G:H4'	45:DW:43:THR:H	1.36	0.88
5:AE:151:LEU:HD13	8:AH:77:GLU:HG2	1.56	0.88
29:BG:89:ILE:HG12	29:BG:162:ILE:HG22	1.56	0.88
23:BA:197:A:H8	23:BA:197:A:H5'	1.36	0.88
23:DA:197:A:H5'	23:DA:197:A:H8	1.36	0.88
43:BU:8:LYS:HZ2	43:BU:8:LYS:H	1.19	0.87
47:DY:17:SER:HB3	47:DY:18:PRO:HD3	1.57	0.87
29:BG:101:ARG:H	29:BG:101:ARG:HE	1.18	0.87
23:BA:857:C:H4'	45:BW:23:VAL:HG21	1.55	0.87
33:BK:3:GLN:HB2	33:BK:4:PRO:HD2	1.54	0.87
50:B2:35:GLU:HB2	50:B2:49:CYS:SG	2.14	0.87
1:CA:922:G:H4'	5:CE:20:GLN:HA	1.56	0.87

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
41:DS:24:ILE:HG21	41:DS:36:LEU:HD21	1.56	0.87
23:DA:142:G:H4'	42:DT:35:THR:HG21	1.55	0.87
27:BE:67:GLN:O	27:BE:67:GLN:HG3	1.73	0.87
1:CA:93:U:H2'	1:CA:95:G:C8	2.09	0.87
50:D2:35:GLU:HB2	50:D2:49:CYS:SG	2.15	0.87
23:BA:270(J):G:HO2'	23:BA:270(K):G:H8	1.19	0.86
21:AU:22:ARG:HD2	21:AU:23:PRO:HD2	1.56	0.86
9:CI:103:THR:HG22	9:CI:105:ASP:H	1.40	0.86
43:BU:81:LYS:HD2	43:BU:96:ILE:HD12	1.57	0.86
23:BA:2426:A:H3'	23:BA:2427:C:H5''	1.56	0.86
1:AA:1223:C:H5'	1:AA:1224:G:H5''	1.55	0.86
49:B1:50:THR:HG22	49:B1:51:TYR:H	1.38	0.86
9:AI:103:THR:HG22	9:AI:105:ASP:H	1.40	0.86
23:BA:676:A:H8	23:BA:2069:G:H21	1.24	0.86
23:BA:954:G:H5''	35:BM:13:GLN:HG3	1.57	0.86
23:DA:857:C:H4'	45:DW:23:VAL:HG21	1.55	0.86
23:DA:1899:G:H21	23:DA:1902:C:N4	1.73	0.86
23:BA:2331:G:H4'	45:BW:43:THR:H	1.38	0.86
10:AJ:48:THR:HA	10:AJ:62:HIS:HB3	1.55	0.86
1:AA:1261:A:H62	1:AA:1274:G:H21	1.19	0.86
27:DE:67:GLN:O	27:DE:67:GLN:HG3	1.74	0.85
1:AA:93:U:H2'	1:AA:95:G:C8	2.11	0.85
3:CC:105:GLU:HG2	3:CC:106:VAL:H	1.40	0.85
37:BO:35:ILE:HG12	37:BO:101:LEU:HD21	1.57	0.85
39:DQ:88:ILE:HB	39:DQ:90:VAL:HG12	1.58	0.85
23:BA:2681:C:H5	23:BA:2725:A:N6	1.74	0.85
47:BY:17:SER:HB3	47:BY:18:PRO:HD3	1.56	0.85
23:BA:2712:U:H1'	23:BA:712(B):A:C8	2.11	0.85
1:AA:979:C:H3'	1:AA:980:C:H5''	1.56	0.85
49:B1:59:VAL:HG12	49:B1:60:GLU:H	1.40	0.85
43:BU:88:LYS:HE2	43:BU:93:GLY:HA3	1.58	0.85
42:BT:35:THR:O	42:BT:39:ILE:HG12	1.77	0.85
37:DO:35:ILE:HG12	37:DO:101:LEU:HD21	1.57	0.85
33:DK:119:PRO:HB2	38:DP:68:TYR:HE1	1.41	0.85
49:D1:59:VAL:HG12	49:D1:60:GLU:H	1.40	0.85
23:DA:2681:C:H5	23:DA:2725:A:N6	1.74	0.85
42:DT:35:THR:O	42:DT:39:ILE:HG12	1.76	0.85
5:CE:151:LEU:HD13	8:CH:77:GLU:HG2	1.56	0.85
39:BQ:88:ILE:HB	39:BQ:90:VAL:HG12	1.59	0.85
10:CJ:48:THR:HA	10:CJ:62:HIS:HB3	1.55	0.85
30:DH:5:LEU:HD23	30:DH:5:LEU:H	1.40	0.85

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
43:DU:2:ARG:HG2	43:DU:3:VAL:HG23	1.59	0.85
1:AA:91:C:H2'	1:AA:92:G:H8	1.42	0.85
3:AC:105:GLU:HG2	3:AC:106:VAL:H	1.41	0.85
17:CQ:9:VAL:HG12	17:CQ:56:VAL:HG22	1.58	0.84
28:DF:41:GLN:HG2	28:DF:155:MET:HB3	1.59	0.84
23:BA:2439:A:C8	23:BA:2439:A:H5'	2.13	0.84
30:BH:5:LEU:HD23	30:BH:5:LEU:H	1.39	0.84
43:BU:45:VAL:HG22	43:BU:62:GLU:HB3	1.59	0.84
23:DA:2712:U:H1'	23:DA:712(B):A:C8	2.12	0.84
23:BA:1813:G:H1'	25:BC:50:THR:HG21	1.59	0.84
43:BU:2:ARG:HG2	43:BU:3:VAL:HG23	1.59	0.84
34:DL:41:ARG:HH22	34:DL:45:LEU:HB2	1.43	0.84
33:BK:119:PRO:HB2	38:BP:68:TYR:HE1	1.39	0.84
28:BF:41:GLN:HG2	28:BF:155:MET:HB3	1.60	0.84
23:DA:2426:A:H3'	23:DA:2427:C:C5'	2.08	0.84
1:CA:1346:A:H61	1:CA:1374:A:H3'	1.42	0.84
46:DX:11:ARG:HB3	46:DX:12:PRO:HD2	1.60	0.84
23:DA:2744:G:H21	29:DG:143:GLN:HE22	1.23	0.84
34:DL:45:LEU:HD23	34:DL:46:LYS:H	1.43	0.84
46:BX:11:ARG:HB3	46:BX:12:PRO:HD2	1.59	0.84
17:AQ:9:VAL:HG12	17:AQ:56:VAL:HG22	1.57	0.84
1:CA:91:C:H2'	1:CA:92:G:H8	1.43	0.83
23:DA:245:G:H5'	34:DL:73:GLY:HA2	1.59	0.83
29:DG:89:ILE:HG12	29:DG:162:ILE:HG22	1.58	0.83
5:CE:51:VAL:HB	5:CE:52:PRO:HD3	1.59	0.83
23:DA:270(J):G:HO2'	23:DA:270(K):G:H8	1.24	0.83
45:DW:35:ASN:H	45:DW:35:ASN:HD22	1.25	0.83
23:BA:2426:A:H3'	23:BA:2427:C:C5'	2.08	0.83
28:DF:55:LYS:HD2	28:DF:58:GLN:HE21	1.42	0.83
26:DD:91:VAL:HB	26:DD:95:ILE:HD11	1.60	0.83
38:DP:26:ASP:HB2	38:DP:91:ARG:HA	1.61	0.83
23:DA:848:G:H2'	23:DA:849:A:C8	2.13	0.83
41:DS:13:SER:HB3	41:DS:16:LYS:HD2	1.59	0.83
23:DA:2439:A:H5'	23:DA:2439:A:C8	2.14	0.83
25:BC:133:LEU:HD23	25:BC:136:ILE:HD12	1.60	0.83
44:DV:97:GLU:HB3	44:DV:125:LEU:HD21	1.61	0.83
36:BN:54:LEU:HD11	36:BN:65:LEU:HD23	1.61	0.83
25:BC:159:ALA:HB1	25:BC:198:ASN:O	1.78	0.83
23:BA:848:G:H2'	23:BA:849:A:C8	2.14	0.83
6:AF:99:ALA:HB2	18:AR:31:LEU:HD22	1.61	0.83
43:DU:45:VAL:HG22	43:DU:62:GLU:HB3	1.59	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
44:BV:97:GLU:HB3	44:BV:125:LEU:HD21	1.60	0.82
23:BA:1899:G:H21	23:BA:1902:C:N4	1.74	0.82
1:AA:1346:A:H61	1:AA:1374:A:H3'	1.42	0.82
1:AA:942:G:H21	9:AI:124:GLN:HE22	1.26	0.82
38:BP:26:ASP:HB2	38:BP:91:ARG:HA	1.60	0.82
25:DC:159:ALA:HB1	25:DC:198:ASN:O	1.78	0.82
28:BF:55:LYS:HD2	28:BF:58:GLN:HE21	1.42	0.82
43:DU:8:LYS:N	43:DU:8:LYS:HZ2	1.77	0.82
23:BA:2744:G:H21	29:BG:143:GLN:HE22	1.25	0.82
23:DA:676:A:H8	23:DA:2069:G:H21	1.25	0.82
10:CJ:75:ILE:HG13	10:CJ:76:ASN:H	1.44	0.82
11:CK:22:HIS:HB3	11:CK:29:ILE:HG13	1.61	0.82
1:CA:105:G:H2'	1:CA:106:C:C6	2.15	0.82
34:DL:89:ALA:HB1	34:DL:121:LYS:HD3	1.62	0.82
13:CM:99:ARG:HB2	13:CM:101:GLN:HE21	1.45	0.82
23:DA:2327:A:H2'	23:DA:2328:A:C8	2.15	0.82
33:BK:119:PRO:HB2	38:BP:68:TYR:CE1	2.14	0.82
1:AA:673:G:H2'	1:AA:674:G:C8	2.15	0.82
29:DG:92:ILE:H	29:DG:92:ILE:HD12	1.45	0.82
11:AK:22:HIS:HB3	11:AK:29:ILE:HG13	1.60	0.81
45:BW:35:ASN:HD22	45:BW:35:ASN:H	1.24	0.81
52:B4:8:ASN:ND2	52:B4:11:LYS:H	1.78	0.81
17:CQ:12:SER:HB3	17:CQ:20:THR:HB	1.62	0.81
34:BL:41:ARG:HH22	34:BL:45:LEU:HB2	1.44	0.81
25:DC:133:LEU:HD23	25:DC:136:ILE:HD12	1.61	0.81
25:DC:25:THR:O	25:DC:27:THR:HG22	1.80	0.81
23:DA:2400:G:H4'	51:D3:19:ARG:HD3	1.61	0.81
43:DU:88:LYS:HE2	43:DU:93:GLY:HA3	1.62	0.81
27:BE:41:LEU:HA	27:BE:44:ARG:HD3	1.62	0.81
23:BA:947:G:C2'	23:BA:948:G:H5''	2.10	0.81
23:BA:245:G:H5'	34:BL:73:GLY:HA2	1.61	0.81
23:DA:1899:G:N2	23:DA:1902:C:H41	1.79	0.81
43:DU:78:ALA:HB3	43:DU:81:LYS:HE3	1.62	0.81
43:BU:78:ALA:HB3	43:BU:81:LYS:HE3	1.63	0.81
10:AJ:75:ILE:HG13	10:AJ:76:ASN:H	1.43	0.81
26:BD:91:VAL:HB	26:BD:95:ILE:HD11	1.61	0.81
13:AM:99:ARG:HB2	13:AM:101:GLN:HE21	1.46	0.81
23:BA:2327:A:H2'	23:BA:2328:A:C8	2.15	0.81
34:BL:89:ALA:HB1	34:BL:121:LYS:HD3	1.62	0.81
1:CA:1129:C:H4'	1:CA:1130:A:H5''	1.62	0.81
36:DN:54:LEU:HD11	36:DN:65:LEU:HD23	1.60	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
53:D5:54:GLU:HA	53:D5:57:ARG:HH12	1.46	0.81
1:AA:690:G:H2'	1:AA:691:G:C8	2.15	0.81
43:DU:81:LYS:HD3	43:DU:97:ARG:HB3	1.63	0.81
35:DM:81:VAL:HG12	35:DM:82:ARG:HG2	1.62	0.81
29:DG:162:ILE:H	29:DG:162:ILE:HD13	1.45	0.81
1:CA:673:G:H2'	1:CA:674:G:C8	2.15	0.81
3:AC:195:VAL:HG12	3:AC:196:LEU:H	1.46	0.81
23:BA:2400:G:H4'	51:B3:19:ARG:HD3	1.61	0.81
28:DF:109:VAL:HG11	28:DF:142:PRO:HG3	1.62	0.81
41:BS:13:SER:HB3	41:BS:16:LYS:HD2	1.62	0.81
5:AE:78:HIS:HE1	5:AE:143:ARG:H	1.27	0.81
29:BG:162:ILE:H	29:BG:162:ILE:HD13	1.46	0.81
30:DH:71:ILE:HG23	30:DH:72:LEU:HD22	1.63	0.81
23:DA:1813:G:H1'	25:DC:50:THR:HG21	1.63	0.81
1:AA:1348:U:H4'	9:AI:120:ARG:HD2	1.63	0.80
1:CA:1347:G:C8	9:CI:107:ARG:HB3	2.15	0.80
39:BQ:83:LEU:HG	39:BQ:88:ILE:HD11	1.63	0.80
26:BD:117:MET:HE1	26:BD:136:ARG:HA	1.62	0.80
1:AA:82:U:H2'	1:AA:85:U:H5	1.46	0.80
3:CC:195:VAL:HG12	3:CC:196:LEU:H	1.46	0.80
23:DA:947:G:H2'	23:DA:948:G:H5''	1.63	0.80
5:CE:78:HIS:HE1	5:CE:143:ARG:H	1.27	0.80
1:AA:38:G:H22	1:AA:397:A:H5'	1.47	0.80
1:AA:105:G:H2'	1:AA:106:C:C6	2.17	0.80
1:CA:82:U:H2'	1:CA:85:U:H5	1.46	0.80
27:DE:41:LEU:HA	27:DE:44:ARG:HD3	1.62	0.80
23:DA:1658:C:OP1	26:DD:132:HIS:ND1	2.15	0.80
6:CF:99:ALA:HB2	18:CR:31:LEU:HD22	1.62	0.80
23:DA:947:G:C2'	23:DA:948:G:H5''	2.11	0.80
1:CA:942:G:H21	9:CI:124:GLN:HE22	1.26	0.80
1:CA:690:G:H2'	1:CA:691:G:C8	2.16	0.80
43:BU:81:LYS:HD3	43:BU:97:ARG:HB3	1.63	0.80
33:DK:119:PRO:HB2	38:DP:68:TYR:CE1	2.16	0.80
13:CM:60:VAL:HG13	13:CM:64:TRP:HE1	1.47	0.80
23:BA:2873:A:C2	36:BN:6:SER:HB2	2.17	0.80
28:BF:109:VAL:HG11	28:BF:142:PRO:HG3	1.61	0.79
1:AA:1346:A:H5'	9:AI:120:ARG:HH12	1.45	0.79
30:DH:62:LYS:HB2	30:DH:133:HIS:CE1	2.18	0.79
23:BA:996:A:H4'	39:BQ:92:ARG:HH12	1.45	0.79
35:BM:81:VAL:HG12	35:BM:82:ARG:HG2	1.63	0.79
23:DA:2873:A:C2	36:DN:6:SER:HB2	2.17	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:CC:141:VAL:HG11	3:CC:202:ILE:HD12	1.63	0.79
3:CC:43:LEU:O	3:CC:47:LEU:HB3	1.82	0.79
34:BL:45:LEU:HD23	34:BL:46:LYS:H	1.44	0.79
42:DT:29:TRP:CZ3	42:DT:78:LYS:HG3	2.16	0.79
23:DA:2025:C:H2'	23:DA:2026:C:C6	2.18	0.79
29:BG:92:ILE:HD12	29:BG:92:ILE:H	1.47	0.79
52:D4:8:ASN:ND2	52:D4:11:LYS:H	1.79	0.79
27:BE:65:TRP:CZ3	27:BE:75:HIS:HD2	2.01	0.79
48:BZ:8:LEU:HD12	48:BZ:31:LEU:HA	1.63	0.79
23:BA:1579:A:H5'	23:BA:1579:A:H8	1.45	0.79
34:BL:6:LEU:H	34:BL:6:LEU:HD23	1.48	0.79
23:DA:948:G:H8	23:DA:948:G:H5'	1.47	0.79
3:AC:141:VAL:HG11	3:AC:202:ILE:HD12	1.65	0.79
1:CA:1443:G:H3'	1:CA:1446:A:C5'	2.11	0.79
1:CA:1224:G:H4'	13:CM:102:ARG:HH22	1.48	0.79
44:DV:77:ASP:HB2	44:DV:84:GLU:HG3	1.63	0.79
1:AA:1443:G:H3'	1:AA:1446:A:C5'	2.11	0.79
1:AA:957:U:H4'	19:AS:79:THR:HB	1.63	0.79
23:BA:773:U:C4'	25:BC:47:GLY:HA3	2.13	0.79
1:CA:957:U:H4'	19:CS:79:THR:HB	1.64	0.79
34:DL:6:LEU:HD23	34:DL:6:LEU:H	1.46	0.79
5:AE:43:LEU:HD11	5:AE:132:ALA:HB1	1.65	0.79
25:DC:264:LYS:O	25:DC:267:SER:HB2	1.83	0.79
45:DW:35:ASN:ND2	45:DW:35:ASN:H	1.81	0.79
25:BC:25:THR:O	25:BC:27:THR:HG22	1.83	0.79
23:BA:2637:U:H5''	26:BD:82:ARG:NH2	1.97	0.79
1:AA:1129:C:H4'	1:AA:1130:A:H5''	1.62	0.78
23:BA:2394:C:OP1	34:BL:63:PRO:HD2	1.83	0.78
2:AB:101:MET:HA	2:AB:108:ILE:HG13	1.65	0.78
33:BK:35:VAL:HG11	33:BK:103:ALA:HB3	1.65	0.78
1:AA:977:A:H2'	1:AA:978:A:H5''	1.65	0.78
2:CB:101:MET:HA	2:CB:108:ILE:HG13	1.65	0.78
44:BV:77:ASP:HB2	44:BV:84:GLU:HG3	1.63	0.78
1:CA:977:A:H2'	1:CA:978:A:H5''	1.65	0.78
23:DA:996:A:H4'	39:DQ:92:ARG:HH12	1.45	0.78
23:DA:1021:A:H62	23:DA:1141:U:H3	1.31	0.78
39:DQ:92:ARG:NH1	39:DQ:92:ARG:HB2	1.97	0.78
35:DM:75:THR:HA	35:DM:88:GLY:HA2	1.64	0.78
5:CE:43:LEU:HD11	5:CE:132:ALA:HB1	1.66	0.78
2:AB:77:ALA:HB2	2:AB:211:ILE:HD13	1.64	0.78
27:DE:65:TRP:CZ3	27:DE:75:HIS:HD2	2.01	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DA:125:G:H4'	23:DA:126:A:OP2	1.82	0.78
34:DL:36:LYS:HG3	34:DL:41:ARG:HB2	1.63	0.78
53:B5:54:GLU:HA	53:B5:57:ARG:HH12	1.47	0.78
28:DF:66:GLN:HG2	28:DF:67:LYS:H	1.48	0.78
16:AP:4:ILE:HG13	16:AP:21:VAL:HG12	1.65	0.78
37:DO:49:VAL:HG12	37:DO:73:LEU:HD23	1.65	0.78
23:BA:947:G:H2'	23:BA:948:G:H5''	1.64	0.78
3:AC:150:LYS:HB3	3:AC:201:TYR:HB2	1.66	0.78
25:BC:71:ASP:HB3	25:BC:103:ARG:HH22	1.49	0.78
47:DY:16:LEU:HB2	47:DY:20:GLU:HG3	1.63	0.78
23:BA:2025:C:H2'	23:BA:2026:C:C6	2.17	0.78
1:CA:1279:A:H62	3:CC:26:LYS:HE2	1.49	0.78
47:BY:16:LEU:HB2	47:BY:20:GLU:HG3	1.64	0.78
37:DO:24:LEU:O	37:DO:86:ALA:HB3	1.82	0.78
25:BC:244:ARG:HG3	25:BC:245:PRO:HD2	1.65	0.78
48:DZ:8:LEU:HD12	48:DZ:31:LEU:HA	1.66	0.78
37:BO:24:LEU:O	37:BO:86:ALA:HB3	1.82	0.78
1:CA:922:G:H2'	1:CA:923:A:C8	2.19	0.78
1:CA:91:C:H2'	1:CA:92:G:C8	2.19	0.78
30:BH:62:LYS:HB2	30:BH:133:HIS:CE1	2.19	0.78
30:BH:71:ILE:HG23	30:BH:72:LEU:HD22	1.64	0.78
1:CA:38:G:H22	1:CA:397:A:H5'	1.48	0.78
23:BA:1311:G:H5'	23:BA:1311:G:H8	1.48	0.78
23:DA:773:U:C4'	25:DC:47:GLY:HA3	2.14	0.78
13:AM:60:VAL:HG13	13:AM:64:TRP:HE1	1.47	0.78
42:DT:63:LYS:NZ	42:DT:72:LYS:HB3	1.99	0.78
1:AA:673:G:H5''	6:AF:87:ARG:NH1	2.00	0.77
45:BW:35:ASN:H	45:BW:35:ASN:ND2	1.81	0.77
42:BT:29:TRP:CZ3	42:BT:78:LYS:HG3	2.19	0.77
1:AA:328:C:H4'	1:AA:329:A:H5'	1.65	0.77
23:DA:1006:C:O2	32:DJ:129:MET:HG2	1.83	0.77
23:DA:587:C:N3	34:DL:33:ARG:HD2	2.00	0.77
23:BA:1021:A:H62	23:BA:1141:U:H3	1.32	0.77
23:BA:1899:G:N2	23:BA:1902:C:H41	1.79	0.77
35:BM:8:LYS:HG3	35:BM:9:TYR:H	1.50	0.77
34:DL:122:PRO:HA	34:DL:141:ALA:O	1.84	0.77
23:DA:2210:G:N3	23:DA:2210:G:H3'	1.99	0.77
23:BA:587:C:N3	34:BL:33:ARG:HD2	2.00	0.77
35:BM:75:THR:HA	35:BM:88:GLY:HA2	1.65	0.77
3:CC:150:LYS:HB3	3:CC:201:TYR:HB2	1.65	0.77
23:DA:1311:G:H5'	23:DA:1311:G:H8	1.48	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1348:U:H4'	9:CI:120:ARG:HD2	1.65	0.77
1:AA:1391:U:H2'	1:AA:1392:G:C8	2.20	0.77
25:BC:264:LYS:O	25:BC:267:SER:HB2	1.84	0.77
23:BA:1006:C:O2	32:BJ:129:MET:HG2	1.84	0.77
25:DC:33:LEU:HD23	25:DC:33:LEU:H	1.50	0.77
23:BA:1658:C:OP1	26:BD:132:HIS:ND1	2.16	0.77
8:AH:42:GLU:HG3	8:AH:109:ILE:HD12	1.66	0.77
23:DA:1579:A:H5'	23:DA:1579:A:H8	1.47	0.77
34:BL:36:LYS:HG3	34:BL:41:ARG:HB2	1.67	0.77
25:BC:242:ARG:HD3	25:BC:242:ARG:N	1.98	0.77
27:DE:53:THR:HG23	27:DE:55:GLY:H	1.49	0.77
1:AA:922:G:H2'	1:AA:923:A:C8	2.20	0.77
17:AQ:12:SER:HB3	17:AQ:20:THR:HB	1.65	0.77
41:DS:1:MET:HE2	41:DS:2:GLU:H	1.50	0.77
39:DQ:83:LEU:HG	39:DQ:88:ILE:HD11	1.65	0.77
28:BF:66:GLN:HG2	28:BF:67:LYS:H	1.48	0.77
35:DM:75:THR:HA	35:DM:88:GLY:CA	2.15	0.77
1:AA:91:C:H2'	1:AA:92:G:C8	2.19	0.77
38:BP:27:THR:HG23	38:BP:89:VAL:HG13	1.67	0.77
23:BA:1210:A:H5''	23:BA:1210:A:C8	2.20	0.77
25:DC:30:GLU:CD	25:DC:63:ARG:HE	1.87	0.77
23:BA:1210:A:H5''	23:BA:1210:A:H8	1.48	0.77
23:DA:1175:U:H2'	23:DA:1176:G:H8	1.50	0.77
39:BQ:24:TYR:HB2	39:BQ:29:SER:HB3	1.65	0.77
39:BQ:92:ARG:HB2	39:BQ:92:ARG:NH1	1.98	0.77
23:BA:773:U:H4'	25:BC:47:GLY:HA3	1.66	0.77
7:AG:69:VAL:HA	7:AG:138:LYS:HD2	1.67	0.77
39:DQ:24:TYR:HB2	39:DQ:29:SER:HB3	1.65	0.77
25:DC:71:ASP:HB3	25:DC:103:ARG:HH22	1.49	0.77
1:CA:1346:A:H5'	9:CI:120:ARG:HH12	1.50	0.76
38:DP:27:THR:HG23	38:DP:89:VAL:HG13	1.66	0.76
23:BA:125:G:H4'	23:BA:126:A:OP2	1.84	0.76
23:DA:2394:C:OP1	34:DL:63:PRO:HD2	1.85	0.76
37:BO:49:VAL:HG12	37:BO:73:LEU:HD23	1.67	0.76
52:D4:12:ARG:NH2	52:D4:44:PRO:HB3	2.00	0.76
23:DA:2637:U:H5''	26:DD:82:ARG:NH2	2.00	0.76
12:AL:82:VAL:HG23	12:AL:106:ALA:HB2	1.68	0.76
37:DO:103:GLU:O	37:DO:107:GLU:HG2	1.84	0.76
2:CB:20:GLU:HB2	2:CB:190:THR:HB	1.67	0.76
43:DU:27:VAL:HG12	43:DU:39:VAL:HG22	1.68	0.76
27:BE:53:THR:HG23	27:BE:55:GLY:H	1.50	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:DD:201:THR:HG22	26:DD:202:LYS:H	1.50	0.76
23:BA:2210:G:N3	23:BA:2210:G:H3'	2.00	0.76
10:AJ:92:THR:HG23	10:AJ:93:GLY:H	1.50	0.76
1:CA:1327:C:OP1	21:CU:20:LYS:HB3	1.85	0.76
40:DR:4:ILE:HB	40:DR:39:LEU:HB2	1.68	0.76
23:BA:1510:A:H2'	23:BA:1511:A:C8	2.21	0.76
1:CA:328:C:H4'	1:CA:329:A:H5'	1.68	0.76
1:CA:1014:A:H5'	19:CS:14:HIS:CD2	2.21	0.76
47:BY:39:ALA:HA	47:BY:45:SER:HB3	1.66	0.76
1:CA:1056:U:H5'	3:CC:163:ALA:HB2	1.68	0.76
25:BC:33:LEU:H	25:BC:33:LEU:HD23	1.50	0.76
16:CP:4:ILE:HG13	16:CP:21:VAL:HG12	1.65	0.76
30:BH:82:ARG:HB3	30:BH:89:TYR:HB2	1.67	0.76
1:AA:1378:C:H5''	7:AG:6:ARG:HE	1.50	0.76
23:BA:2377:A:H2'	23:BA:2378:A:C8	2.21	0.76
23:BA:948:G:H5'	23:BA:948:G:H8	1.48	0.76
23:DA:1510:A:H2'	23:DA:1511:A:C8	2.21	0.76
3:AC:43:LEU:O	3:AC:47:LEU:HB3	1.84	0.76
13:AM:67:GLU:HG3	13:AM:68:GLY:H	1.50	0.76
26:BD:201:THR:HG22	26:BD:202:LYS:H	1.50	0.76
12:CL:82:VAL:HG23	12:CL:106:ALA:HB2	1.68	0.76
23:DA:140:A:H8	23:DA:1408:C:HO2'	1.34	0.76
3:CC:14:ILE:HG23	3:CC:15:THR:H	1.51	0.76
23:BA:140:A:H8	23:BA:1408:C:HO2'	1.31	0.76
43:DU:31:LEU:HD23	43:DU:31:LEU:H	1.50	0.76
1:CA:134:A:H61	16:CP:25:ARG:NH1	1.83	0.76
23:DA:1210:A:H8	23:DA:1210:A:H5''	1.51	0.76
34:BL:115:LEU:HA	34:BL:134:ALA:HB2	1.67	0.75
23:DA:773:U:H4'	25:DC:47:GLY:HA3	1.68	0.75
47:DY:39:ALA:HA	47:DY:45:SER:HB3	1.67	0.75
2:AB:91:PRO:HA	2:AB:154:LEU:HD11	1.68	0.75
23:DA:2377:A:H2'	23:DA:2378:A:C8	2.20	0.75
34:DL:115:LEU:HA	34:DL:134:ALA:HB2	1.69	0.75
26:BD:2:LYS:HD3	26:BD:95:ILE:HG22	1.67	0.75
23:BA:1175:U:H2'	23:BA:1176:G:H8	1.50	0.75
8:CH:42:GLU:HG3	8:CH:109:ILE:HD12	1.66	0.75
2:CB:77:ALA:HB2	2:CB:211:ILE:HD13	1.67	0.75
37:BO:103:GLU:O	37:BO:107:GLU:HG2	1.86	0.75
34:DL:45:LEU:HD23	34:DL:46:LYS:N	2.01	0.75
23:DA:2542:A:N3	23:DA:2542:A:H5''	2.01	0.75
26:DD:52:LEU:HD12	26:DD:52:LEU:H	1.50	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
29:BG:35:VAL:HG21	29:BG:75:ALA:HB2	1.66	0.75
23:BA:603:A:N1	23:BA:655:A:H1'	2.00	0.75
42:BT:63:LYS:NZ	42:BT:72:LYS:HB3	2.01	0.75
34:BL:122:PRO:HA	34:BL:141:ALA:O	1.86	0.75
52:D4:8:ASN:C	52:D4:8:ASN:HD22	1.90	0.75
26:DD:201:THR:HG22	26:DD:202:LYS:N	2.01	0.75
23:DA:528:A:H8	23:DA:528:A:H3'	1.52	0.75
23:DA:603:A:N1	23:DA:655:A:H1'	2.02	0.75
32:DJ:66:THR:H	32:DJ:71:MET:HE3	1.52	0.75
34:BL:45:LEU:HD23	34:BL:46:LYS:N	2.01	0.75
30:BH:83:ALA:HB2	30:BH:88:ILE:HD13	1.69	0.75
36:DN:10:LEU:HB2	36:DN:17:ARG:HE	1.52	0.75
35:BM:75:THR:HA	35:BM:88:GLY:CA	2.16	0.75
2:CB:91:PRO:HA	2:CB:154:LEU:HD11	1.68	0.75
7:CG:69:VAL:HA	7:CG:138:LYS:HD2	1.67	0.75
23:BA:343:C:H5'	23:BA:343:C:H6	1.50	0.75
13:CM:67:GLU:HG3	13:CM:68:GLY:H	1.50	0.75
34:BL:33:ARG:CG	34:BL:36:LYS:HD3	2.13	0.75
23:BA:2427:C:H6	23:BA:2427:C:H5'	1.51	0.75
9:CI:13:ALA:HB2	9:CI:68:GLY:HA3	1.68	0.75
23:BA:2039:C:H2'	23:BA:2040:C:H6	1.52	0.75
26:BD:201:THR:HG22	26:BD:202:LYS:N	2.03	0.74
23:DA:343:C:H5'	23:DA:343:C:H6	1.49	0.74
26:DD:117:MET:HE1	26:DD:136:ARG:HA	1.69	0.74
40:BR:66:ARG:HD2	40:BR:88:ARG:CZ	2.17	0.74
40:BR:4:ILE:HB	40:BR:39:LEU:HB2	1.67	0.74
19:AS:50:ALA:HB1	19:AS:57:HIS:HB3	1.70	0.74
28:BF:86:MET:SD	28:BF:87:PRO:HD3	2.27	0.74
41:BS:1:MET:HE2	41:BS:2:GLU:H	1.51	0.74
32:DJ:57:LEU:O	32:DJ:72:GLY:HA3	1.87	0.74
26:DD:2:LYS:HD3	26:DD:95:ILE:HG22	1.67	0.74
2:AB:20:GLU:HB2	2:AB:190:THR:HB	1.67	0.74
1:AA:1220:G:H21	19:AS:54:GLY:HA2	1.52	0.74
52:B4:11:LYS:O	52:B4:15:THR:HG23	1.87	0.74
28:DF:28:VAL:O	28:DF:31:VAL:HG12	1.88	0.74
23:DA:2039:C:H2'	23:DA:2040:C:H6	1.52	0.74
25:BC:227:ASN:HB3	25:BC:228:PRO:HD2	1.69	0.74
25:BC:30:GLU:CD	25:BC:63:ARG:HE	1.90	0.74
25:DC:244:ARG:HG3	25:DC:245:PRO:HD2	1.68	0.74
43:BU:27:VAL:HG12	43:BU:39:VAL:HG22	1.69	0.74
23:DA:2056:G:N2	50:D2:4:HIS:O	2.20	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DA:547:A:H2'	23:DA:548:A:H8	1.51	0.74
1:CA:1371:G:OP1	9:CI:11:LYS:HB3	1.87	0.74
23:DA:674:G:H1'	27:DE:74:ARG:HD3	1.69	0.74
43:BU:71:LYS:NZ	43:BU:71:LYS:HB2	2.03	0.74
8:CH:10:LEU:HD22	8:CH:83:ILE:HD11	1.69	0.74
23:DA:2427:C:H6	23:DA:2427:C:H5'	1.52	0.74
36:DN:10:LEU:HB2	36:DN:17:ARG:NE	2.03	0.74
3:AC:30:ARG:HD3	14:AN:38:GLY:HA3	1.70	0.74
42:BT:50:LYS:H	42:BT:87:GLN:HE22	1.35	0.74
3:CC:30:ARG:HD3	14:CN:38:GLY:HA3	1.70	0.74
42:BT:30:VAL:HG11	42:BT:39:ILE:HD12	1.69	0.74
1:AA:199:G:H2'	1:AA:200:G:H5''	1.70	0.74
10:AJ:74:ILE:HD13	10:AJ:74:ILE:H	1.53	0.74
33:DK:71:ARG:HH12	38:DP:74:ARG:HH22	1.32	0.74
33:BK:71:ARG:HH12	38:BP:74:ARG:HH22	1.36	0.74
23:BA:1019:U:H3	23:BA:114(B):A:H62	1.36	0.74
35:DM:8:LYS:HG3	35:DM:9:TYR:H	1.51	0.74
42:DT:30:VAL:HG11	42:DT:39:ILE:HD12	1.69	0.74
23:DA:2287:A:O2'	23:DA:2288:A:H5''	1.88	0.74
24:BB:13:A:N7	24:BB:70:C:H4'	2.02	0.74
42:DT:15:GLU:H	42:DT:15:GLU:CD	1.91	0.74
25:BC:118:VAL:HG22	25:BC:119:ALA:H	1.53	0.74
10:CJ:92:THR:HG23	10:CJ:93:GLY:H	1.50	0.74
30:DH:83:ALA:HB2	30:DH:88:ILE:HD13	1.68	0.74
38:DP:84:GLN:HG3	38:DP:85:LYS:HG3	1.70	0.74
23:BA:2056:G:N2	50:B2:4:HIS:O	2.21	0.73
32:BJ:157:ARG:N	32:BJ:158:PRO:HD3	2.03	0.73
44:BV:48:PHE:HA	44:BV:51:ALA:HB3	1.70	0.73
34:DL:26:GLY:HA2	34:DL:30:THR:HG23	1.69	0.73
34:DL:33:ARG:CG	34:DL:36:LYS:HD3	2.13	0.73
32:DJ:157:ARG:H	32:DJ:158:PRO:CD	2.01	0.73
23:BA:2025:C:H2'	23:BA:2026:C:H6	1.51	0.73
42:DT:50:LYS:H	42:DT:87:GLN:HE22	1.34	0.73
23:BA:2542:A:H5''	23:BA:2542:A:N3	2.02	0.73
23:DA:1314:C:H6	23:DA:1314:C:H5'	1.53	0.73
13:CM:9:ILE:HG22	13:CM:11:ARG:HG3	1.71	0.73
23:BA:528:A:H3'	23:BA:528:A:H8	1.53	0.73
24:BB:43:C:H4'	28:BF:98:ARG:HH12	1.52	0.73
32:BJ:157:ARG:H	32:BJ:158:PRO:CD	2.01	0.73
32:DJ:157:ARG:N	32:DJ:158:PRO:HD3	2.03	0.73
30:DH:82:ARG:HB3	30:DH:89:TYR:HB2	1.70	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:AC:14:ILE:HG23	3:AC:15:THR:H	1.51	0.73
38:BP:84:GLN:HG3	38:BP:85:LYS:HG3	1.70	0.73
5:CE:39:GLY:HA2	5:CE:69:VAL:HB	1.70	0.73
23:DA:2025:C:H2'	23:DA:2026:C:H6	1.53	0.73
9:AI:13:ALA:HB2	9:AI:68:GLY:HA3	1.68	0.73
25:BC:186:HIS:HD2	25:BC:188:GLU:H	1.37	0.73
32:BJ:57:LEU:O	32:BJ:72:GLY:HA3	1.88	0.73
24:DB:43:C:H4'	28:DF:98:ARG:HH12	1.52	0.73
40:BR:5:VAL:HG23	40:BR:37:VAL:HG23	1.70	0.73
28:DF:86:MET:SD	28:DF:87:PRO:HD3	2.29	0.73
2:AB:84:GLU:HB3	2:AB:219:VAL:HG21	1.71	0.73
44:DV:126:VAL:HG12	44:DV:163:LEU:HA	1.71	0.73
12:AL:74:HIS:CD2	12:AL:76:LEU:H	2.07	0.73
23:DA:528:A:H2	23:DA:2043:C:H4'	1.52	0.73
24:DB:13:A:N7	24:DB:70:C:H4'	2.04	0.73
25:DC:147:LEU:HD13	25:DC:155:LEU:HD11	1.69	0.73
26:DD:51:PHE:HD1	26:DD:52:LEU:HG	1.53	0.73
30:BH:9:LEU:HB3	30:BH:12:LEU:HD23	1.71	0.73
37:BO:30:ARG:HB3	37:BO:35:ILE:HD13	1.71	0.73
44:BV:126:VAL:HG12	44:BV:163:LEU:HA	1.71	0.73
28:DF:77:ILE:HG22	28:DF:80:PHE:H	1.54	0.73
36:BN:10:LEU:HB2	36:BN:17:ARG:HE	1.53	0.73
23:BA:2287:A:O2'	23:BA:2288:A:H5''	1.88	0.73
23:BA:806:C:OP2	34:BL:39:LYS:HD2	1.89	0.73
1:CA:736:C:H2'	1:CA:737:A:H8	1.53	0.73
23:BA:2469:A:H2	23:BA:2481:G:H21	1.35	0.73
23:BA:973:A:OP2	40:BR:78:LYS:NZ	2.21	0.73
2:CB:60:ASP:O	2:CB:64:ARG:HG2	1.89	0.73
19:CS:50:ALA:HB1	19:CS:57:HIS:HB3	1.69	0.73
29:DG:35:VAL:HG21	29:DG:75:ALA:HB2	1.70	0.73
42:BT:15:GLU:H	42:BT:15:GLU:CD	1.91	0.73
23:DA:1678:G:O5'	23:DA:1678:G:H8	1.72	0.73
43:BU:96:ILE:HD11	43:BU:99:CYS:HB2	1.69	0.72
37:DO:30:ARG:HB3	37:DO:35:ILE:HD13	1.72	0.72
53:B5:57:ARG:HB2	53:B5:57:ARG:NH1	2.04	0.72
1:CA:1123:A:H4'	10:CJ:36:GLY:HA3	1.71	0.72
23:DA:2469:A:H2	23:DA:2481:G:H21	1.34	0.72
34:DL:29:LYS:HD2	34:DL:29:LYS:N	2.04	0.72
10:CJ:96:ILE:HD13	10:CJ:96:ILE:H	1.54	0.72
1:CA:1129:C:H4'	1:CA:1130:A:C5'	2.19	0.72
36:BN:10:LEU:HB2	36:BN:17:ARG:NE	2.04	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
28:BF:77:ILE:HG22	28:BF:80:PHE:H	1.53	0.72
27:BE:139:PHE:HB2	27:BE:166:ALA:HB1	1.71	0.72
19:AS:6:LYS:HG2	19:AS:7:LYS:HD3	1.71	0.72
43:DU:71:LYS:HB2	43:DU:71:LYS:NZ	2.04	0.72
35:BM:75:THR:HG21	35:BM:85:LYS:NZ	2.04	0.72
1:AA:716:A:H1'	11:AK:118:GLY:HA2	1.69	0.72
26:DD:154:LYS:HA	26:DD:154:LYS:HE3	1.71	0.72
43:DU:96:ILE:HD11	43:DU:99:CYS:HB2	1.70	0.72
23:BA:547:A:H2'	23:BA:548:A:H8	1.51	0.72
25:DC:30:GLU:HG3	25:DC:63:ARG:NH2	2.04	0.72
23:DA:1210:A:H5''	23:DA:1210:A:C8	2.24	0.72
17:CQ:69:LYS:C	17:CQ:70:ARG:HD2	2.10	0.72
5:AE:39:GLY:HA2	5:AE:69:VAL:HB	1.71	0.72
28:BF:28:VAL:O	28:BF:31:VAL:HG12	1.89	0.72
49:B1:38:ALA:HA	49:B1:55:PRO:HA	1.72	0.72
23:DA:848:G:C4	23:DA:933:A:H8	2.07	0.72
44:DV:48:PHE:HA	44:DV:51:ALA:HB3	1.71	0.72
25:BC:30:GLU:HG3	25:BC:63:ARG:NH2	2.04	0.72
23:BA:1314:C:H5'	23:BA:1314:C:H6	1.55	0.72
23:DA:1405:U:H2'	23:DA:1406:U:C6	2.25	0.72
5:CE:78:HIS:CE1	5:CE:143:ARG:H	2.08	0.72
23:BA:848:G:C4	23:BA:933:A:H8	2.07	0.72
53:D5:57:ARG:HB2	53:D5:57:ARG:NH1	2.05	0.72
34:DL:59:LEU:HA	34:DL:61:ARG:NE	2.04	0.72
46:DX:17:SER:HB3	46:DX:44:PRO:HD3	1.70	0.72
2:AB:55:PHE:HE1	2:AB:218:ALA:HA	1.54	0.72
10:AJ:96:ILE:H	10:AJ:96:ILE:HD13	1.54	0.72
16:CP:27:LYS:H	16:CP:27:LYS:HD2	1.55	0.72
16:AP:27:LYS:HD2	16:AP:27:LYS:H	1.53	0.72
5:CE:6:PHE:HD2	5:CE:36:ASP:HB3	1.54	0.72
40:DR:66:ARG:HD2	40:DR:88:ARG:CZ	2.20	0.72
26:BD:52:LEU:HD12	26:BD:52:LEU:H	1.53	0.72
26:BD:51:PHE:HD1	26:BD:52:LEU:HG	1.53	0.72
25:DC:242:ARG:HD3	25:DC:242:ARG:N	2.00	0.72
34:BL:64:LYS:HB2	53:B5:25:MET:HG3	1.72	0.72
47:DY:2:LYS:CD	47:DY:2:LYS:H	2.03	0.72
23:BA:106:C:H1'	43:BU:2:ARG:HE	1.54	0.72
4:AD:49:ARG:NH2	4:AD:50:ARG:HB2	2.04	0.72
1:AA:1320:C:H42	19:AS:36:ARG:HG3	1.52	0.72
33:DK:35:VAL:HG11	33:DK:103:ALA:HB3	1.69	0.72
1:CA:1378:C:H5''	7:CG:6:ARG:NE	2.05	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DA:106:C:H1'	43:DU:2:ARG:HE	1.52	0.72
48:BZ:43:ILE:O	48:BZ:47:VAL:HG23	1.89	0.72
26:BD:179:GLU:HB3	26:BD:181:LEU:HD23	1.72	0.72
2:CB:55:PHE:HE1	2:CB:218:ALA:HA	1.54	0.72
15:CO:63:ARG:NH2	15:CO:87:ILE:HG21	2.03	0.72
23:BA:1021:A:C3'	23:BA:1021:A:C8	2.72	0.72
12:CL:74:HIS:CD2	12:CL:76:LEU:H	2.07	0.72
23:BA:528:A:H2	23:BA:2043:C:H4'	1.54	0.72
23:BA:2287:A:H62	23:BA:2344:U:H3	1.37	0.72
23:BA:556:G:H2'	23:BA:557:U:C6	2.24	0.72
27:DE:160:ASN:OD1	27:DE:163:VAL:HG23	1.90	0.72
32:BJ:65:TRP:HA	32:BJ:71:MET:HE1	1.72	0.72
30:BH:110:ASP:HB2	30:BH:113:ARG:HG2	1.70	0.72
1:AA:1422:G:H5''	33:BK:48:PRO:HB3	1.71	0.72
2:CB:84:GLU:HB3	2:CB:219:VAL:HG21	1.71	0.72
42:BT:63:LYS:HD2	42:BT:72:LYS:HA	1.72	0.71
52:B4:35:ARG:HG3	52:B4:42:LEU:HD11	1.71	0.71
11:CK:21:ILE:HG13	11:CK:30:VAL:HG12	1.71	0.71
27:BE:155:LEU:HD23	27:BE:186:ILE:HD13	1.71	0.71
35:DM:43:THR:HB	35:DM:45:GLN:HE21	1.55	0.71
23:DA:1019:U:H3	23:DA:114(B):A:H62	1.37	0.71
10:AJ:50:ILE:HB	14:AN:41:ARG:NH2	2.05	0.71
44:BV:163:LEU:HD23	44:BV:163:LEU:H	1.55	0.71
25:BC:27:THR:O	25:BC:27:THR:HG23	1.91	0.71
11:AK:21:ILE:HG13	11:AK:30:VAL:HG12	1.71	0.71
26:DD:179:GLU:HB3	26:DD:181:LEU:HD23	1.72	0.71
23:DA:1516:U:H2'	23:DA:1517:G:H8	1.55	0.71
19:CS:18:LYS:HG2	19:CS:31:ILE:HD13	1.72	0.71
6:AF:23:LYS:O	6:AF:27:GLN:HG2	1.88	0.71
6:CF:16:GLN:H	6:CF:16:GLN:CD	1.93	0.71
23:BA:2543:G:H8	23:BA:2543:G:H5'	1.53	0.71
23:DA:2075:U:H2'	23:DA:2238:G:N2	2.06	0.71
23:BA:2075:U:H2'	23:BA:2238:G:N2	2.05	0.71
51:B3:42:TRP:HA	51:B3:42:TRP:CE3	2.25	0.71
53:D5:26:LYS:HA	53:D5:48:PHE:HE2	1.55	0.71
23:DA:1429:G:H2'	23:DA:1430:C:C6	2.25	0.71
1:CA:392:G:H2'	1:CA:393:A:H8	1.54	0.71
5:AE:6:PHE:HD2	5:AE:36:ASP:HB3	1.54	0.71
23:BA:1343:G:H5'	23:BA:1343:G:C8	2.24	0.71
34:BL:29:LYS:N	34:BL:29:LYS:HD2	2.05	0.71
1:CA:1220:G:H21	19:CS:54:GLY:HA2	1.55	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:BL:49:ARG:CG	34:BL:50:ARG:N	2.52	0.71
52:B4:8:ASN:HD22	52:B4:8:ASN:C	1.92	0.71
23:BA:1311:G:C8	23:BA:1311:G:H5'	2.26	0.71
42:DT:63:LYS:HD2	42:DT:72:LYS:HA	1.71	0.71
17:AQ:69:LYS:C	17:AQ:70:ARG:HD2	2.10	0.71
25:BC:147:LEU:HD13	25:BC:155:LEU:HD11	1.71	0.71
1:CA:199:G:H2'	1:CA:200:G:H5''	1.71	0.71
27:DE:139:PHE:HB2	27:DE:166:ALA:HB1	1.73	0.71
23:DA:2415:G:H4'	34:DL:66:GLY:CA	2.20	0.71
10:CJ:50:ILE:HB	14:CN:41:ARG:NH2	2.04	0.71
25:DC:35:LYS:HG3	25:DC:104:TYR:CE2	2.26	0.71
34:BL:59:LEU:HA	34:BL:61:ARG:NE	2.05	0.71
8:AH:10:LEU:HD22	8:AH:83:ILE:HD11	1.71	0.71
31:DI:14:LYS:HE2	31:DI:14:LYS:HA	1.71	0.71
4:CD:49:ARG:NH2	4:CD:50:ARG:HB2	2.04	0.71
23:BA:1429:G:H2'	23:BA:1430:C:C6	2.26	0.71
1:AA:736:C:H2'	1:AA:737:A:H8	1.54	0.71
12:CL:74:HIS:HD2	12:CL:76:LEU:H	1.38	0.71
23:DA:1336:A:H2'	23:DA:1337:G:H8	1.55	0.71
48:DZ:43:ILE:O	48:DZ:47:VAL:HG23	1.91	0.71
46:BX:17:SER:HB3	46:BX:44:PRO:HD3	1.69	0.71
23:BA:1516:U:H2'	23:BA:1517:G:H8	1.56	0.71
1:AA:1129:C:H4'	1:AA:1130:A:C5'	2.19	0.71
28:BF:83:ARG:HG3	28:BF:84:LYS:H	1.56	0.71
53:B5:26:LYS:HA	53:B5:48:PHE:HE2	1.55	0.71
34:BL:16:ARG:HE	34:BL:17:LYS:N	1.89	0.71
19:CS:6:LYS:HG2	19:CS:7:LYS:HD3	1.71	0.71
23:DA:556:G:H2'	23:DA:557:U:C6	2.26	0.71
23:BA:2415:G:H4'	34:BL:66:GLY:CA	2.20	0.71
37:BO:34:HIS:HA	37:BO:54:LEU:HD23	1.72	0.71
2:AB:54:THR:HG21	2:AB:201:ILE:HD11	1.72	0.71
2:CB:54:THR:HG21	2:CB:201:ILE:HD11	1.72	0.71
28:DF:83:ARG:HG3	28:DF:84:LYS:H	1.54	0.71
37:DO:34:HIS:HA	37:DO:54:LEU:HD23	1.72	0.71
10:CJ:74:ILE:H	10:CJ:74:ILE:HD13	1.54	0.71
23:BA:1336:A:H2'	23:BA:1337:G:H8	1.55	0.71
40:DR:5:VAL:HG23	40:DR:37:VAL:HG23	1.71	0.71
23:DA:197:A:H5'	23:DA:197:A:C8	2.24	0.71
34:DL:16:ARG:HE	34:DL:17:LYS:N	1.88	0.71
28:BF:19:LEU:HD11	28:BF:172:LEU:HD13	1.72	0.71
23:DA:1343:G:H5'	23:DA:1343:G:C8	2.26	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DM:75:THR:HG21	35:DM:85:LYS:NZ	2.06	0.71
13:AM:9:ILE:HG22	13:AM:11:ARG:HG3	1.71	0.71
18:AR:50:ILE:HD12	18:AR:70:ILE:HG21	1.73	0.71
25:DC:227:ASN:HB3	25:DC:228:PRO:HD2	1.72	0.71
30:DH:9:LEU:HB3	30:DH:12:LEU:HD23	1.73	0.71
4:AD:4:TYR:CE1	4:AD:11:LEU:HD11	2.26	0.71
35:BM:43:THR:HB	35:BM:45:GLN:HE21	1.54	0.71
43:BU:31:LEU:H	43:BU:31:LEU:HD23	1.54	0.71
4:CD:20:TYR:HD2	4:CD:26:CYS:HB3	1.56	0.71
23:DA:1541:U:H3'	23:DA:1542:G:O3'	1.91	0.70
7:CG:15:ASP:HB3	7:CG:19:GLY:H	1.56	0.70
39:DQ:15:LYS:O	39:DQ:19:LYS:HG3	1.91	0.70
46:DX:50:ARG:HG2	46:DX:59:THR:HG22	1.73	0.70
1:AA:168:G:H2'	1:AA:169:C:H5''	1.73	0.70
23:DA:1021:A:C8	23:DA:1021:A:C3'	2.73	0.70
23:DA:2392:A:H2	23:DA:2424:C:H42	1.39	0.70
38:BP:26:ASP:CB	38:BP:91:ARG:HA	2.21	0.70
1:CA:82:U:H2'	1:CA:85:U:C5	2.25	0.70
7:AG:69:VAL:HG22	7:AG:135:VAL:HG22	1.72	0.70
25:DC:155:LEU:HD23	25:DC:177:LEU:HD21	1.72	0.70
6:AF:16:GLN:H	6:AF:16:GLN:CD	1.95	0.70
6:CF:23:LYS:O	6:CF:27:GLN:HG2	1.91	0.70
25:DC:118:VAL:HG22	25:DC:119:ALA:H	1.54	0.70
11:CK:110:ASP:HB3	18:CR:85:LEU:HB3	1.74	0.70
5:AE:78:HIS:CE1	5:AE:143:ARG:H	2.08	0.70
25:DC:27:THR:O	25:DC:27:THR:HG23	1.91	0.70
52:D4:11:LYS:O	52:D4:15:THR:HG23	1.89	0.70
23:DA:322:A:H3'	27:DE:169:ASN:ND2	2.06	0.70
28:BF:7:LEU:HD23	28:BF:10:LYS:HD2	1.74	0.70
23:BA:796:C:H2'	23:BA:797:C:C6	2.25	0.70
23:DA:860:U:H5	23:DA:917:A:N7	1.89	0.70
23:DA:806:C:OP2	34:DL:39:LYS:HD2	1.90	0.70
7:CG:69:VAL:HG22	7:CG:135:VAL:HG22	1.72	0.70
23:BA:2740:A:H2'	23:BA:2741:A:C8	2.26	0.70
28:DF:19:LEU:HD11	28:DF:172:LEU:HD13	1.73	0.70
22:CV:6189:A:H2'	22:CV:6190:C:H6	1.56	0.70
44:DV:108:PRO:HG3	44:DV:141:VAL:HG22	1.73	0.70
30:DH:76:THR:HG22	30:DH:141:LYS:HD3	1.74	0.70
25:DC:144:ALA:HB3	25:DC:192:THR:HG23	1.73	0.70
7:CG:102:ARG:HG2	7:CG:106:GLN:HE21	1.56	0.70
23:BA:141(A):A:H8	23:BA:1595:G:H21	1.38	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
44:DV:104:PHE:HA	44:DV:139:VAL:HB	1.73	0.70
19:AS:18:LYS:HG2	19:AS:31:ILE:HD13	1.72	0.70
32:DJ:90:LEU:HD12	32:DJ:90:LEU:H	1.54	0.70
2:AB:60:ASP:O	2:AB:64:ARG:HG2	1.91	0.70
23:BA:674:G:H1'	27:BE:74:ARG:HD3	1.73	0.70
23:DA:2036:C:H6	23:DA:2036:C:H5'	1.56	0.70
49:D1:38:ALA:HA	49:D1:55:PRO:HA	1.71	0.70
23:DA:1899:G:N2	23:DA:1902:C:N4	2.38	0.70
30:BH:56:LYS:HA	30:BH:59:ALA:HB3	1.73	0.70
51:D3:15:GLU:OE2	51:D3:18:ARG:HD2	1.92	0.70
23:BA:2261:C:C6	45:BW:16:SER:HB3	2.27	0.70
23:DA:2261:C:C6	45:DW:16:SER:HB3	2.25	0.70
52:B4:12:ARG:NH2	52:B4:44:PRO:HB3	2.05	0.70
23:DA:1657:C:H2'	23:DA:1658:C:H6	1.54	0.70
6:CF:69:GLU:O	6:CF:72:VAL:HG12	1.92	0.70
30:BH:76:THR:HG22	30:BH:141:LYS:HD3	1.74	0.70
34:BL:26:GLY:HA2	34:BL:30:THR:HG23	1.71	0.70
23:BA:2036:C:H5'	23:BA:2036:C:H6	1.57	0.70
52:B4:5:TRP:NE1	52:B4:7:PRO:HG3	2.06	0.70
23:BA:2068:U:N3	23:BA:2430:A:H2	1.89	0.70
43:BU:8:LYS:N	43:BU:8:LYS:HZ2	1.90	0.70
44:BV:76:LEU:H	44:BV:76:LEU:HD12	1.56	0.70
47:DY:16:LEU:O	47:DY:20:GLU:HB2	1.92	0.70
23:DA:2543:G:H8	23:DA:2543:G:H5'	1.57	0.70
23:DA:528:A:C8	23:DA:528:A:H3'	2.26	0.70
25:BC:201:HIS:O	25:BC:204:ILE:HG13	1.90	0.70
37:BO:90:GLY:O	37:BO:92:TYR:N	2.25	0.70
1:AA:392:G:H2'	1:AA:393:A:H8	1.56	0.70
1:CA:729:A:H2'	1:CA:730:G:H8	1.57	0.70
3:AC:6:HIS:HD2	3:AC:7:PRO:HD2	1.56	0.70
3:CC:6:HIS:HD2	3:CC:7:PRO:HD2	1.56	0.70
3:AC:206:GLU:HG2	3:AC:207:VAL:HG23	1.73	0.70
27:DE:155:LEU:HD23	27:DE:186:ILE:HD13	1.72	0.70
28:BF:121:ASN:HD22	28:BF:122:PRO:HD2	1.56	0.70
23:DA:2740:A:H2'	23:DA:2741:A:C8	2.26	0.70
14:AN:24:CYS:HB3	14:AN:29:ARG:H	1.57	0.70
32:BJ:90:LEU:H	32:BJ:90:LEU:HD12	1.56	0.70
31:BI:14:LYS:HA	31:BI:14:LYS:HE2	1.72	0.70
34:DL:49:ARG:CG	34:DL:50:ARG:N	2.53	0.69
42:BT:27:THR:HB	42:BT:80:ILE:HB	1.74	0.69
47:BY:2:LYS:CD	47:BY:2:LYS:H	2.05	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
42:BT:50:LYS:H	42:BT:87:GLN:NE2	1.90	0.69
46:DX:27:GLU:HB2	46:DX:33:LYS:HA	1.74	0.69
8:AH:86:ILE:HB	8:AH:133:LEU:HD22	1.74	0.69
4:AD:20:TYR:HD2	4:AD:26:CYS:HB3	1.56	0.69
37:DO:90:GLY:O	37:DO:92:TYR:N	2.25	0.69
22:AV:6189:A:H2'	22:AV:6190:C:H6	1.56	0.69
23:BA:760:G:H2'	23:BA:761:A:H5'	1.74	0.69
48:DZ:8:LEU:CD1	48:DZ:31:LEU:HD12	2.22	0.69
23:BA:1541:U:H3'	23:BA:1542:G:O3'	1.92	0.69
25:DC:186:HIS:HD2	25:DC:188:GLU:H	1.38	0.69
52:D4:5:TRP:NE1	52:D4:7:PRO:HG3	2.07	0.69
29:DG:43:VAL:HA	29:DG:52:VAL:HG22	1.73	0.69
39:BQ:15:LYS:O	39:BQ:19:LYS:HG3	1.92	0.69
23:DA:1418:G:O5'	23:DA:1418:G:H8	1.74	0.69
29:DG:17:VAL:HG22	29:DG:26:VAL:HG22	1.74	0.69
15:AO:63:ARG:NH2	15:AO:87:ILE:HG21	2.04	0.69
42:DT:27:THR:HB	42:DT:80:ILE:HB	1.74	0.69
34:DL:64:LYS:HB2	53:D5:25:MET:HG3	1.73	0.69
23:DA:1311:G:H5'	23:DA:1311:G:C8	2.27	0.69
12:AL:74:HIS:HD2	12:AL:76:LEU:H	1.38	0.69
11:AK:87:THR:HA	11:AK:91:ARG:HH21	1.57	0.69
23:BA:1678:G:O5'	23:BA:1678:G:H8	1.75	0.69
25:BC:35:LYS:HG3	25:BC:104:TYR:CE2	2.26	0.69
1:AA:922:G:H4'	5:AE:20:GLN:HA	1.74	0.69
1:AA:522:C:H41	12:AL:52:ARG:HH22	1.39	0.69
28:DF:121:ASN:HD22	28:DF:122:PRO:HD2	1.56	0.69
4:CD:4:TYR:CE1	4:CD:11:LEU:HD11	2.26	0.69
11:CK:18:ARG:HB3	11:CK:33:THR:HG23	1.73	0.69
30:DH:110:ASP:HB2	30:DH:113:ARG:HG2	1.72	0.69
6:AF:7:ASN:HD21	18:AR:34:TYR:HE1	1.40	0.69
1:CA:1224:G:H4'	13:CM:102:ARG:NH2	2.07	0.69
44:DV:163:LEU:H	44:DV:163:LEU:HD23	1.56	0.69
1:CA:1152:A:OP1	10:CJ:68:HIS:HD2	1.75	0.69
13:AM:14:ARG:NH1	13:AM:42:ALA:HA	2.06	0.69
23:BA:2023:G:H5'	23:BA:2617:C:H4'	1.74	0.69
3:CC:206:GLU:HG2	3:CC:207:VAL:HG23	1.73	0.69
1:AA:82:U:H2'	1:AA:85:U:C5	2.25	0.69
23:DA:1174:A:H3'	23:DA:1175:U:H5''	1.73	0.69
46:BX:83:GLU:HG2	46:BX:84:GLY:H	1.57	0.69
39:BQ:50:ARG:HH22	40:BR:72:VAL:HG12	1.58	0.69
32:BJ:36:TRP:HB2	32:BJ:156:GLN:HB2	1.75	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
11:CK:87:THR:HA	11:CK:91:ARG:HH21	1.58	0.69
18:CR:50:ILE:HD12	18:CR:70:ILE:HG21	1.75	0.69
30:DH:56:LYS:HA	30:DH:59:ALA:HB3	1.74	0.69
4:CD:189:PRO:HB2	4:CD:194:LEU:HD21	1.75	0.69
26:BD:154:LYS:HA	26:BD:154:LYS:HE3	1.74	0.69
23:DA:796:C:H2'	23:DA:797:C:C6	2.27	0.69
46:BX:50:ARG:HG2	46:BX:59:THR:HG22	1.73	0.69
23:BA:1163:G:C2'	23:BA:1164:G:H5''	2.23	0.69
23:DA:1046:A:H3'	23:DA:1047:G:H5''	1.75	0.69
23:BA:1174:A:H3'	23:BA:1175:U:H5''	1.73	0.69
2:CB:187:LEU:HA	2:CB:201:ILE:HB	1.74	0.69
23:BA:2415:G:H4'	34:BL:66:GLY:HA3	1.75	0.69
23:BA:1542:G:H1'	23:BA:1543:A:C4	2.28	0.69
30:DH:78:THR:HA	30:DH:143:SER:HB3	1.75	0.69
1:CA:1427:U:H2'	1:CA:1428:A:C8	2.28	0.69
1:CA:168:G:H2'	1:CA:169:C:H5''	1.73	0.69
51:D3:42:TRP:CE3	51:D3:42:TRP:HA	2.25	0.69
35:BM:60:ARG:H	44:BV:179:ASP:HB2	1.56	0.69
46:DX:83:GLU:HG2	46:DX:84:GLY:H	1.57	0.69
1:CA:1226:C:H2'	13:CM:103:THR:HB	1.75	0.69
50:D2:45:VAL:HG12	50:D2:46:CYS:H	1.58	0.69
23:DA:2287:A:H62	23:DA:2344:U:H3	1.38	0.69
30:DH:79:ILE:HB	30:DH:144:VAL:HA	1.73	0.69
23:BA:1544:C:H3'	23:BA:1545:A:H5'	1.75	0.69
27:BE:160:ASN:OD1	27:BE:163:VAL:HG23	1.93	0.69
11:AK:110:ASP:HB3	18:AR:85:LEU:HB3	1.74	0.69
6:AF:47:ARG:HH12	6:AF:56:PRO:HB2	1.58	0.69
23:BA:1899:G:N2	23:BA:1902:C:N4	2.39	0.69
23:BA:1046:A:H1'	31:BI:4:LYS:CD	2.23	0.69
38:DP:26:ASP:CB	38:DP:91:ARG:HA	2.22	0.69
23:DA:2415:G:H4'	34:DL:66:GLY:HA3	1.75	0.69
23:DA:1778:U:H2'	23:DA:1784:A:N6	2.07	0.69
3:AC:19:GLU:HG2	3:AC:40:ARG:HH22	1.56	0.69
23:BA:2062:A:O2'	23:BA:2063:C:H5'	1.93	0.69
25:DC:238:GLY:O	25:DC:239:ARG:C	2.32	0.68
51:B3:42:TRP:HA	51:B3:42:TRP:HE3	1.58	0.68
51:B3:15:GLU:OE2	51:B3:18:ARG:HD2	1.92	0.68
35:DM:60:ARG:H	44:DV:179:ASP:HB2	1.57	0.68
50:B2:45:VAL:HG12	50:B2:46:CYS:H	1.56	0.68
8:CH:86:ILE:HB	8:CH:133:LEU:HD22	1.73	0.68
46:BX:11:ARG:HB3	46:BX:12:PRO:CD	2.21	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:DC:155:LEU:HD23	25:DC:177:LEU:CD2	2.23	0.68
23:DA:1516:U:H2'	23:DA:1517:G:C8	2.28	0.68
44:BV:104:PHE:HA	44:BV:139:VAL:HB	1.74	0.68
3:CC:95:THR:HG22	3:CC:96:GLY:H	1.58	0.68
39:BQ:55:ARG:HA	39:BQ:58:ARG:HD2	1.73	0.68
46:BX:27:GLU:HB2	46:BX:33:LYS:HA	1.75	0.68
35:DM:24:GLY:HA2	35:DM:101:ARG:HA	1.75	0.68
1:AA:1435:G:H2'	1:AA:1436:U:C6	2.28	0.68
1:AA:729:A:H2'	1:AA:730:G:H8	1.59	0.68
39:DQ:90:VAL:HG13	39:DQ:91:ASP:H	1.58	0.68
1:CA:939:G:H5''	7:CG:102:ARG:HH12	1.58	0.68
30:BH:79:ILE:HB	30:BH:144:VAL:HA	1.75	0.68
23:BA:591:C:O2	53:B5:2:PRO:HA	1.94	0.68
28:DF:7:LEU:HD23	28:DF:10:LYS:HD2	1.74	0.68
14:CN:24:CYS:HB3	14:CN:29:ARG:H	1.56	0.68
1:AA:1286:A:C8	21:AU:22:ARG:NH2	2.61	0.68
46:DX:11:ARG:HB3	46:DX:12:PRO:CD	2.23	0.68
34:DL:59:LEU:HA	34:DL:61:ARG:CZ	2.24	0.68
1:CA:673:G:H5''	6:CF:87:ARG:NH1	2.08	0.68
23:BA:1516:U:H2'	23:BA:1517:G:C8	2.29	0.68
24:DB:79:C:H2'	24:DB:80:U:O4'	1.94	0.68
1:CA:1369:C:H2'	1:CA:1370:G:C8	2.28	0.68
35:BM:24:GLY:HA2	35:BM:101:ARG:HA	1.75	0.68
12:CL:56:LYS:HG2	12:CL:66:THR:HG22	1.75	0.68
23:BA:2056:G:H2'	23:BA:2056:G:N3	2.08	0.68
23:BA:733:G:N7	23:BA:761:A:C6	2.62	0.68
7:AG:102:ARG:HG2	7:AG:106:GLN:HE21	1.57	0.68
23:BA:528:A:C8	23:BA:528:A:H3'	2.28	0.68
23:DA:1542:G:H1'	23:DA:1543:A:C4	2.28	0.68
12:AL:56:LYS:HG2	12:AL:66:THR:HG22	1.74	0.68
23:BA:587:C:N4	34:BL:33:ARG:HG2	2.08	0.68
23:DA:2068:U:N3	23:DA:2430:A:H2	1.90	0.68
42:DT:50:LYS:H	42:DT:87:GLN:NE2	1.90	0.68
1:CA:1512:U:H2'	1:CA:1513:A:C8	2.29	0.68
1:AA:1056:U:H5'	3:AC:163:ALA:HB2	1.74	0.68
23:BA:860:U:H5	23:BA:917:A:N7	1.91	0.68
3:AC:31:HIS:O	3:AC:35:GLU:HG2	1.94	0.68
19:AS:69:HIS:HB3	19:AS:73:GLU:HG3	1.76	0.68
4:AD:3:ARG:HH21	4:AD:118:ARG:HD3	1.58	0.68
24:BB:55:U:H4'	28:BF:27:ASN:HD21	1.58	0.68
23:DA:2023:G:H5'	23:DA:2617:C:H4'	1.75	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DA:1826:G:H4'	25:DC:242:ARG:HE	1.59	0.68
3:CC:195:VAL:HG12	3:CC:196:LEU:N	2.08	0.68
25:BC:155:LEU:HD23	25:BC:177:LEU:HD21	1.74	0.68
4:AD:189:PRO:HB2	4:AD:194:LEU:HD21	1.75	0.68
51:D3:36:LEU:HD23	51:D3:36:LEU:H	1.58	0.68
29:BG:43:VAL:HA	29:BG:52:VAL:HG22	1.74	0.68
25:BC:244:ARG:HG3	25:BC:245:PRO:CD	2.24	0.68
23:BA:1316:U:H2'	23:BA:1317:A:C8	2.29	0.68
25:DC:201:HIS:O	25:DC:204:ILE:HG13	1.94	0.68
7:AG:15:ASP:HB3	7:AG:19:GLY:H	1.57	0.68
4:CD:23:GLY:HA3	4:CD:112:VAL:HG22	1.76	0.68
15:AO:5:LYS:N	15:AO:5:LYS:HD3	2.09	0.68
15:CO:5:LYS:N	15:CO:5:LYS:HD3	2.08	0.68
34:BL:146:VAL:HG22	34:BL:147:LEU:H	1.59	0.68
1:AA:1346:A:N1	1:AA:1374:A:H5''	2.09	0.68
25:BC:96:HIS:HD2	25:BC:102:LYS:HG2	1.59	0.68
52:D4:35:ARG:HG3	52:D4:42:LEU:HD11	1.76	0.68
25:DC:148:GLU:HB2	25:DC:151:LYS:HD2	1.75	0.68
51:B3:36:LEU:HD23	51:B3:36:LEU:H	1.59	0.68
23:DA:591:C:O2	53:D5:2:PRO:HA	1.93	0.68
39:BQ:90:VAL:HG13	39:BQ:91:ASP:H	1.58	0.68
4:CD:9:CYS:HB3	4:CD:32:ALA:HB2	1.76	0.68
18:CR:66:LEU:O	18:CR:70:ILE:HG12	1.94	0.68
51:D3:42:TRP:HE3	51:D3:42:TRP:HA	1.58	0.68
6:CF:7:ASN:HD21	18:CR:34:TYR:HE1	1.41	0.68
1:AA:1014:A:H5'	19:AS:14:HIS:CD2	2.28	0.68
43:BU:42:VAL:HG12	43:BU:65:ALA:HB3	1.76	0.68
32:BJ:66:THR:HB	32:BJ:69:VAL:HG12	1.76	0.68
23:DA:973:A:OP2	40:DR:78:LYS:NZ	2.23	0.68
3:CC:31:HIS:O	3:CC:35:GLU:HG2	1.94	0.68
41:DS:18:ARG:HG2	41:DS:76:VAL:CG1	2.24	0.68
1:AA:1369:C:H2'	1:AA:1370:G:C8	2.29	0.68
1:AA:939:G:H5''	7:AG:102:ARG:NH1	2.09	0.67
1:CA:1347:G:H22	1:CA:1373:G:H2'	1.58	0.67
23:DA:673:C:H5''	27:DE:81:PRO:HD2	1.75	0.67
23:BA:1826:G:H4'	25:BC:242:ARG:HE	1.58	0.67
10:CJ:32:ALA:H	10:CJ:78:ASN:HD21	1.43	0.67
2:AB:187:LEU:HA	2:AB:201:ILE:HB	1.75	0.67
23:BA:1175:U:H2'	23:BA:1176:G:C8	2.29	0.67
23:DA:1544:C:H3'	23:DA:1545:A:H5'	1.75	0.67
30:BH:87:LYS:HA	30:BH:122:GLU:HA	1.76	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:BC:217:ARG:HH11	25:BC:217:ARG:HG2	1.59	0.67
4:CD:3:ARG:HH21	4:CD:118:ARG:HD3	1.58	0.67
30:DH:87:LYS:HA	30:DH:122:GLU:HA	1.76	0.67
11:AK:18:ARG:HB3	11:AK:33:THR:HG23	1.74	0.67
6:CF:47:ARG:HH12	6:CF:56:PRO:HB2	1.58	0.67
25:DC:96:HIS:HD2	25:DC:102:LYS:HG2	1.58	0.67
3:AC:17:ASP:HB2	3:AC:21:ARG:HH22	1.59	0.67
23:DA:141(A):A:H8	23:DA:1595:G:H21	1.40	0.67
18:CR:44:LEU:HD22	18:CR:79:LEU:HD22	1.75	0.67
44:BV:108:PRO:HG3	44:BV:141:VAL:HG22	1.75	0.67
44:DV:136:PHE:C	44:DV:137:ILE:HD12	2.15	0.67
8:CH:51:VAL:HG21	8:CH:60:ARG:HG3	1.76	0.67
1:CA:1285:A:H4'	1:CA:1286:A:O5'	1.95	0.67
23:BA:1405:U:H2'	23:BA:1406:U:C6	2.28	0.67
30:BH:92:VAL:HA	30:BH:96:ASP:OD2	1.95	0.67
3:AC:195:VAL:HG12	3:AC:196:LEU:N	2.08	0.67
23:BA:2637:U:H5''	26:BD:82:ARG:HH21	1.59	0.67
47:DY:46:GLN:O	47:DY:47:ASN:HB2	1.93	0.67
23:DA:185:U:H2'	23:DA:186:G:C8	2.29	0.67
32:BJ:118:PRO:O	32:BJ:121:VAL:HG22	1.94	0.67
25:BC:78:LYS:HD3	25:BC:114:GLY:HA2	1.77	0.67
24:DB:66:A:H61	24:DB:107:U:H2'	1.59	0.67
1:AA:1427:U:H2'	1:AA:1428:A:C8	2.29	0.67
48:BZ:8:LEU:CD1	48:BZ:31:LEU:HD12	2.24	0.67
10:AJ:45:ARG:HB2	10:AJ:65:LEU:HB3	1.76	0.67
27:DE:8:GLN:HA	27:DE:21:ALA:HA	1.77	0.67
19:CS:69:HIS:HB3	19:CS:73:GLU:HG3	1.76	0.67
10:CJ:45:ARG:HB2	10:CJ:65:LEU:HB3	1.76	0.67
23:DA:65:C:H2'	23:DA:66:C:H6	1.60	0.67
23:DA:587:C:N4	34:DL:33:ARG:HG2	2.09	0.67
23:DA:760:G:H2'	23:DA:761:A:H5'	1.77	0.67
28:BF:98:ARG:HD2	28:BF:98:ARG:H	1.60	0.67
4:AD:9:CYS:HB3	4:AD:32:ALA:HB2	1.77	0.67
39:DQ:50:ARG:HH22	40:DR:72:VAL:HG12	1.58	0.67
4:AD:108:LEU:HD23	4:AD:110:PHE:HE2	1.60	0.67
23:DA:2056:G:N3	23:DA:2056:G:H2'	2.08	0.67
23:DA:1046:A:H1'	31:DI:4:LYS:CD	2.23	0.67
23:BA:1046:A:H3'	23:BA:1047:G:H5''	1.75	0.67
23:DA:245:G:H2'	23:DA:246:C:H6	1.59	0.67
26:BD:92:THR:HB	26:BD:94:GLU:HG2	1.77	0.67
18:AR:44:LEU:HD22	18:AR:79:LEU:HD22	1.76	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:CC:19:GLU:HG2	3:CC:40:ARG:HH22	1.56	0.67
23:DA:480:A:OP2	43:DU:46:LYS:HE2	1.95	0.67
29:BG:17:VAL:HG22	29:BG:26:VAL:HG22	1.76	0.67
5:CE:149:GLU:O	5:CE:153:LYS:HB2	1.95	0.67
39:DQ:55:ARG:HA	39:DQ:58:ARG:HD2	1.76	0.67
22:AV:6155:C:H2'	22:AV:6156:A:C8	2.30	0.67
1:AA:1347:G:H22	1:AA:1373:G:H2'	1.59	0.67
1:CA:1443:G:N2	38:DP:119:LYS:HA	2.09	0.67
23:BA:1657:C:H2'	23:BA:1658:C:H6	1.59	0.67
32:DJ:66:THR:HB	32:DJ:69:VAL:HG12	1.76	0.67
23:DA:1316:U:H2'	23:DA:1317:A:C8	2.29	0.67
25:BC:144:ALA:HB3	25:BC:192:THR:HG23	1.75	0.67
24:DB:55:U:H4'	28:DF:27:ASN:HD21	1.59	0.67
3:AC:95:THR:HG22	3:AC:96:GLY:H	1.59	0.67
42:BT:26:TYR:HB3	42:BT:92:LEU:HD13	1.77	0.67
23:BA:1336:A:H2'	23:BA:1337:G:C8	2.30	0.67
1:CA:1151:A:O2'	1:CA:1152:A:H8	1.78	0.67
23:BA:443:A:H2'	27:BE:45:ARG:HH12	1.60	0.67
23:DA:443:A:H2'	27:DE:45:ARG:HH12	1.60	0.67
1:AA:541:G:O2'	4:AD:41:GLY:HA2	1.94	0.67
26:BD:30:PRO:HD3	26:BD:180:ASN:ND2	2.10	0.67
23:BA:1858:G:H1'	23:BA:1884:A:N6	2.10	0.67
34:DL:62:LEU:HD11	53:D5:27:THR:HA	1.76	0.66
23:DA:954:G:H5''	35:DM:13:GLN:CG	2.25	0.66
24:BB:13:A:H5'	45:BW:74:ARG:HH21	1.60	0.66
25:DC:69:ARG:HH21	25:DC:130:ALA:HB2	1.59	0.66
53:D5:39:LYS:O	53:D5:43:GLN:HG2	1.96	0.66
23:DA:1348:G:H2'	23:DA:1349:A:H5''	1.77	0.66
1:AA:1073:U:H2'	1:AA:1074:G:H8	1.60	0.66
7:AG:100:ALA:O	7:AG:104:LEU:HD23	1.95	0.66
23:BA:586:A:H5'	27:BE:89:VAL:HG21	1.77	0.66
34:BL:148:LEU:H	34:BL:148:LEU:HD13	1.60	0.66
23:BA:2822:G:H2'	23:BA:2823:A:H5''	1.77	0.66
27:BE:139:PHE:CB	27:BE:166:ALA:HB1	2.25	0.66
23:DA:1794:U:H2'	23:DA:1795:C:H6	1.59	0.66
1:CA:1079:G:H2'	1:CA:1080:A:C8	2.30	0.66
44:DV:24:LEU:HB2	44:DV:41:LEU:HD23	1.77	0.66
26:DD:30:PRO:HD3	26:DD:180:ASN:ND2	2.10	0.66
32:DJ:118:PRO:O	32:DJ:121:VAL:HG22	1.94	0.66
34:BL:62:LEU:HD11	53:B5:27:THR:HA	1.77	0.66
53:B5:22:VAL:HB	53:B5:54:GLU:HG2	1.77	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BA:330:A:C2	23:BA:1210:A:H2'	2.31	0.66
12:AL:69:ILE:HG13	12:AL:99:ILE:HG21	1.77	0.66
30:BH:82:ARG:HB3	30:BH:89:TYR:CB	2.24	0.66
12:CL:69:ILE:HG13	12:CL:99:ILE:HG21	1.76	0.66
3:CC:17:ASP:HB2	3:CC:21:ARG:HH22	1.60	0.66
25:BC:186:HIS:CD2	25:BC:188:GLU:H	2.14	0.66
1:CA:1435:G:H2'	1:CA:1436:U:C6	2.30	0.66
1:CA:7:G:H21	5:CE:121:LYS:HE2	1.60	0.66
23:BA:65:C:H2'	23:BA:66:C:H6	1.60	0.66
23:BA:1504:C:HO2'	23:BA:1505:C:H6	1.41	0.66
43:DU:8:LYS:H	43:DU:8:LYS:NZ	1.89	0.66
1:CA:1346:A:N1	1:CA:1374:A:H5''	2.10	0.66
46:BX:10:LYS:O	46:BX:11:ARG:HG2	1.96	0.66
1:AA:691:G:C6	11:AK:52:GLY:HA2	2.31	0.66
28:DF:98:ARG:HD2	28:DF:98:ARG:H	1.59	0.66
1:CA:277:C:H5''	17:CQ:68:ARG:NH2	2.11	0.66
44:BV:136:PHE:C	44:BV:137:ILE:HD12	2.15	0.66
23:DA:588:U:H2'	23:DA:589:C:C6	2.30	0.66
1:CA:878:G:H5'	8:CH:89:PRO:HG2	1.77	0.66
13:CM:14:ARG:NH1	13:CM:42:ALA:HA	2.09	0.66
39:BQ:92:ARG:HD2	39:BQ:95:LEU:HG	1.77	0.66
25:BC:94:LEU:HB2	25:BC:104:TYR:HE1	1.60	0.66
47:BY:16:LEU:O	47:BY:20:GLU:HB2	1.95	0.66
23:DA:1175:U:H2'	23:DA:1176:G:C8	2.29	0.66
23:DA:528:A:C2	23:DA:2043:C:H4'	2.31	0.66
18:AR:70:ILE:O	18:AR:74:ARG:HG3	1.96	0.66
7:CG:100:ALA:O	7:CG:104:LEU:HD23	1.94	0.66
3:AC:18:TRP:CD1	14:AN:54:PRO:HA	2.31	0.66
34:BL:111:ARG:HG3	34:BL:128:HIS:CG	2.30	0.66
34:DL:146:VAL:HG22	34:DL:147:LEU:H	1.57	0.66
23:DA:1163:G:C2'	23:DA:1164:G:H5''	2.23	0.66
1:AA:955:U:H1'	1:AA:1227:A:N6	2.08	0.66
23:BA:126:A:OP2	52:B4:19:ARG:HB2	1.95	0.66
22:CV:6172:U:H2'	22:CV:6173:G:H8	1.60	0.66
1:AA:878:G:H5'	8:AH:89:PRO:HG2	1.77	0.66
25:DC:217:ARG:HG2	25:DC:217:ARG:HH11	1.60	0.66
43:DU:42:VAL:HG12	43:DU:65:ALA:HB3	1.78	0.66
22:CV:6155:C:H2'	22:CV:6156:A:C8	2.30	0.66
17:AQ:8:GLY:HA3	17:AQ:23:VAL:HG12	1.78	0.66
41:BS:18:ARG:HG2	41:BS:76:VAL:CG1	2.25	0.66
34:DL:148:LEU:H	34:DL:148:LEU:HD13	1.59	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
42:DT:26:TYR:HB3	42:DT:92:LEU:HD13	1.77	0.66
1:AA:1285:A:H4'	1:AA:1286:A:O5'	1.95	0.66
6:AF:76:ALA:O	6:AF:80:ARG:HG2	1.96	0.66
24:BB:66:A:H61	24:BB:107:U:H2'	1.60	0.66
23:BA:1348:G:H2'	23:BA:1349:A:H5''	1.77	0.66
45:DW:37:LEU:O	45:DW:38:VAL:HG23	1.96	0.66
40:BR:40:LEU:HD23	40:BR:47:VAL:HG23	1.78	0.66
1:AA:939:G:H5''	7:AG:102:ARG:HH12	1.61	0.66
1:AA:1512:U:H2'	1:AA:1513:A:C8	2.31	0.66
27:BE:8:GLN:HA	27:BE:21:ALA:HA	1.77	0.66
42:BT:8:ILE:H	42:BT:8:ILE:HD12	1.60	0.66
23:BA:1614:A:N6	41:BS:87:PRO:HA	2.11	0.66
3:CC:18:TRP:CD1	14:CN:54:PRO:HA	2.31	0.66
23:DA:251:A:C5	23:DA:252:G:H1'	2.31	0.66
1:CA:955:U:H1'	1:CA:1227:A:N6	2.09	0.66
32:DJ:36:TRP:HB2	32:DJ:156:GLN:HB2	1.76	0.66
25:DC:94:LEU:HB2	25:DC:104:TYR:HE1	1.60	0.66
25:BC:148:GLU:HB2	25:BC:151:LYS:HD2	1.76	0.66
29:BG:101:ARG:H	29:BG:101:ARG:NE	1.93	0.66
34:BL:59:LEU:HA	34:BL:61:ARG:CZ	2.26	0.66
50:B2:40:LYS:HE2	50:B2:46:CYS:HB3	1.78	0.66
1:CA:1512:U:H2'	1:CA:1513:A:H8	1.61	0.66
23:BA:774:A:H2	23:BA:787:U:HO2'	1.39	0.66
23:BA:1778:U:H2'	23:BA:1784:A:N6	2.11	0.66
25:BC:238:GLY:O	25:BC:239:ARG:C	2.33	0.65
30:DH:92:VAL:HA	30:DH:96:ASP:OD2	1.96	0.65
25:DC:270:ILE:O	25:DC:271:ILE:HG13	1.96	0.65
23:BA:1418:G:H8	23:BA:1418:G:O5'	1.78	0.65
39:BQ:110:VAL:O	39:BQ:114:LYS:HG2	1.96	0.65
3:AC:59:ARG:HG2	3:AC:64:VAL:HG22	1.78	0.65
25:BC:155:LEU:HD23	25:BC:177:LEU:CD2	2.26	0.65
4:CD:22:LYS:HB2	4:CD:26:CYS:SG	2.35	0.65
1:CA:243:A:H4'	1:CA:244:U:O5'	1.96	0.65
23:DA:2359:C:H2'	23:DA:2360:A:C8	2.31	0.65
5:AE:149:GLU:O	5:AE:153:LYS:HB2	1.96	0.65
33:BK:20:MET:HG2	33:BK:21:CYS:O	1.96	0.65
9:CI:113:LYS:H	9:CI:119:ALA:HA	1.61	0.65
4:AD:23:GLY:HA3	4:AD:112:VAL:HG22	1.76	0.65
1:AA:1356:G:H2'	1:AA:1357:A:C8	2.31	0.65
23:BA:251:A:C5	23:BA:252:G:H1'	2.31	0.65
34:DL:111:ARG:HG3	34:DL:128:HIS:CG	2.30	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BA:2392:A:H2	23:BA:2424:C:H42	1.43	0.65
23:DA:675:A:H4'	27:DE:67:GLN:NE2	2.12	0.65
23:DA:270(G):U:H2'	23:DA:270(H):C:C6	2.31	0.65
26:DD:92:THR:HB	26:DD:94:GLU:HG2	1.78	0.65
23:DA:948:G:C8	23:DA:948:G:H5'	2.31	0.65
1:CA:397:A:H3'	1:CA:397:A:N3	2.11	0.65
3:AC:58:GLU:O	3:AC:64:VAL:HA	1.97	0.65
33:DK:103:ALA:HB1	33:DK:105:GLU:OE1	1.97	0.65
3:CC:83:ARG:O	3:CC:87:LEU:HG	1.96	0.65
23:BA:673:C:H5''	27:BE:81:PRO:HD2	1.78	0.65
15:CO:48:LYS:HE2	15:CO:48:LYS:HA	1.79	0.65
24:BB:74:U:H2'	24:BB:75:G:C8	2.31	0.65
22:AV:6172:U:H2'	22:AV:6173:G:H8	1.61	0.65
23:BA:1694:C:H5''	23:BA:1694:C:H6	1.62	0.65
12:CL:26:LEU:CD1	12:CL:27:LYS:H	2.07	0.65
28:BF:32:PRO:HB2	28:BF:172:LEU:HD22	1.77	0.65
22:AV:6189:A:H2'	22:AV:6190:C:C6	2.32	0.65
23:DA:1332:G:C8	23:DA:1332:G:H5'	2.32	0.65
1:CA:1194:U:H2'	1:CA:1195:C:C6	2.31	0.65
13:CM:16:ASP:HB3	13:CM:34:LEU:HD11	1.78	0.65
25:BC:79:VAL:HG21	25:BC:111:LEU:HD11	1.77	0.65
5:AE:91:LEU:HA	5:AE:120:THR:HG22	1.79	0.65
47:BY:46:GLN:O	47:BY:47:ASN:HB2	1.94	0.65
8:AH:51:VAL:HG21	8:AH:60:ARG:HG3	1.77	0.65
24:DB:74:U:H2'	24:DB:75:G:C8	2.32	0.65
34:BL:41:ARG:NH2	34:BL:45:LEU:HD12	2.11	0.65
34:DL:41:ARG:NH2	34:DL:45:LEU:HD12	2.12	0.65
39:DQ:110:VAL:O	39:DQ:114:LYS:HG2	1.97	0.65
40:DR:49:THR:HB	40:DR:50:PRO:HD2	1.78	0.65
23:BA:675:A:H4'	27:BE:67:GLN:NE2	2.11	0.65
23:BA:910:A:H62	35:BM:12:GLN:HA	1.61	0.65
30:BH:83:ALA:HB3	30:BH:123:LEU:HD12	1.79	0.65
37:BO:99:LYS:O	37:BO:103:GLU:HB2	1.97	0.65
23:DA:1336:A:H2'	23:DA:1337:G:C8	2.30	0.65
18:AR:66:LEU:O	18:AR:70:ILE:HG12	1.95	0.65
28:DF:32:PRO:HB2	28:DF:172:LEU:HD22	1.77	0.65
23:DA:389:G:O6	34:DL:71:VAL:HG23	1.97	0.65
1:CA:1073:U:H2'	1:CA:1074:G:H8	1.61	0.65
27:DE:103:LYS:HA	27:DE:106:ARG:HG3	1.77	0.65
23:BA:2359:C:H2'	23:BA:2360:A:C8	2.32	0.65
23:DA:1858:G:H1'	23:DA:1884:A:N6	2.11	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BA:1332:G:C8	23:BA:1332:G:H5'	2.32	0.65
1:CA:939:G:H5''	7:CG:102:ARG:NH1	2.12	0.65
23:DA:2822:G:H2'	23:DA:2823:A:H5''	1.77	0.65
6:AF:69:GLU:O	6:AF:72:VAL:HG12	1.95	0.65
44:BV:24:LEU:HB2	44:BV:41:LEU:HD23	1.77	0.65
23:DA:733:G:N7	23:DA:761:A:C6	2.64	0.65
23:BA:197:A:C8	23:BA:197:A:H5'	2.24	0.65
10:AJ:32:ALA:H	10:AJ:78:ASN:HD21	1.43	0.65
37:DO:99:LYS:O	37:DO:103:GLU:HB2	1.96	0.65
23:BA:342:G:C2'	23:BA:343:C:H5''	2.27	0.65
22:CV:6189:A:H2'	22:CV:6190:C:C6	2.32	0.65
1:CA:1225:A:H2'	1:CA:1225:A:N3	2.11	0.65
23:DA:1607:C:H4'	23:DA:1608:A:O5'	1.97	0.65
23:DA:2599:G:C8	25:DC:237:GLU:HG3	2.32	0.65
40:DR:22:VAL:HG12	40:DR:23:GLU:N	2.10	0.65
1:CA:105:G:H2'	1:CA:106:C:H6	1.61	0.65
53:D5:22:VAL:HB	53:D5:54:GLU:HG2	1.77	0.65
23:DA:342:G:C2'	23:DA:343:C:H5''	2.26	0.65
23:DA:343:C:C6	23:DA:343:C:H5'	2.32	0.65
20:AT:50:GLU:HB3	20:AT:100:ILE:HD13	1.77	0.65
13:AM:16:ASP:HB3	13:AM:34:LEU:HD11	1.78	0.65
39:BQ:92:ARG:HD3	39:BQ:94:ASN:HB3	1.79	0.65
23:DA:910:A:H62	35:DM:12:GLN:HA	1.62	0.65
6:AF:97:PHE:HD2	18:AR:31:LEU:HD21	1.62	0.65
1:CA:942:G:H21	9:CI:124:GLN:NE2	1.95	0.65
3:CC:58:GLU:O	3:CC:64:VAL:HA	1.96	0.65
4:AD:28:SER:HB3	4:AD:29:PRO:HD2	1.79	0.65
51:B3:11:LEU:HG	51:B3:26:ASN:HB2	1.78	0.65
23:DA:811:U:O2	23:DA:1250:G:H2'	1.97	0.65
1:AA:1225:A:N3	1:AA:1225:A:H2'	2.11	0.65
40:BR:22:VAL:HG12	40:BR:23:GLU:N	2.11	0.65
40:BR:49:THR:HB	40:BR:50:PRO:HD2	1.77	0.65
47:BY:6:VAL:O	47:BY:10:LEU:HG	1.96	0.65
25:DC:244:ARG:HG3	25:DC:245:PRO:CD	2.26	0.65
23:BA:860:U:O2'	23:BA:861:A:H5'	1.96	0.65
25:BC:142:VAL:HG23	25:BC:192:THR:O	1.97	0.65
1:AA:1079:G:H2'	1:AA:1080:A:C8	2.32	0.65
25:DC:10:THR:HG23	25:DC:13:ARG:CB	2.27	0.65
3:AC:83:ARG:O	3:AC:87:LEU:HG	1.96	0.65
20:CT:50:GLU:HB3	20:CT:100:ILE:HD13	1.78	0.65
20:CT:26:ASN:HB2	20:CT:71:THR:HG23	1.78	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BA:2724:C:OP1	26:BD:118:LYS:HE3	1.97	0.65
4:CD:108:LEU:HD23	4:CD:110:PHE:HE2	1.61	0.65
23:DA:330:A:C2	23:DA:1210:A:H2'	2.33	0.64
1:AA:430:A:OP1	4:AD:9:CYS:HB2	1.96	0.64
22:AV:6157:A:H61	22:AV:6172:U:H3	1.42	0.64
30:DH:6:LEU:HD23	30:DH:6:LEU:H	1.63	0.64
5:CE:91:LEU:HA	5:CE:120:THR:HG22	1.78	0.64
23:DA:2637:U:H5''	26:DD:82:ARG:HH21	1.61	0.64
23:BA:140:A:H8	23:BA:1408:C:O2'	1.79	0.64
23:BA:528:A:C2	23:BA:2043:C:H4'	2.32	0.64
23:BA:1343:G:H5'	23:BA:1343:G:H8	1.62	0.64
23:DA:661:C:O3'	34:DL:18:ARG:HG2	1.98	0.64
25:BC:158:ALA:HB3	25:BC:161:THR:HG21	1.79	0.64
10:AJ:55:LYS:O	10:AJ:55:LYS:HD2	1.97	0.64
23:BA:2506:U:H5	23:BA:2583:G:H1	1.45	0.64
1:CA:1298:C:H4'	1:CA:1299:A:C8	2.33	0.64
34:DL:9:ASN:N	34:DL:10:PRO:HD3	2.12	0.64
23:BA:322:A:H3'	27:BE:169:ASN:ND2	2.12	0.64
1:AA:438:G:H4'	1:AA:439:A:OP1	1.97	0.64
23:BA:588:U:H2'	23:BA:589:C:C6	2.32	0.64
46:DX:10:LYS:O	46:DX:11:ARG:HG2	1.97	0.64
1:AA:38:G:N2	1:AA:397:A:H5'	2.13	0.64
1:AA:397:A:N3	1:AA:397:A:H3'	2.12	0.64
26:BD:51:PHE:H	26:BD:75:VAL:HB	1.62	0.64
50:B2:40:LYS:CE	50:B2:46:CYS:HB3	2.27	0.64
23:BA:270(G):U:H2'	23:BA:270(H):C:C6	2.32	0.64
1:CA:1356:G:H2'	1:CA:1357:A:C8	2.31	0.64
2:CB:168:THR:OG1	2:CB:192:SER:HA	1.97	0.64
51:D3:11:LEU:HG	51:D3:26:ASN:HB2	1.78	0.64
29:DG:55:PRO:HG2	29:DG:61:HIS:HD2	1.62	0.64
23:BA:2078:C:H2'	23:BA:2079:U:H6	1.62	0.64
23:BA:1794:U:H2'	23:BA:1795:C:H6	1.62	0.64
23:BA:1022:G:H8	32:BJ:92:GLN:NE2	1.96	0.64
25:DC:131:LEU:CD1	25:DC:136:ILE:HG12	2.28	0.64
1:AA:79:G:H2'	1:AA:80:G:C8	2.32	0.64
6:CF:97:PHE:HD2	18:CR:31:LEU:HD21	1.63	0.64
23:DA:1652:A:OP1	36:DN:9:LYS:HE3	1.98	0.64
19:CS:29:ARG:HB2	19:CS:48:THR:H	1.63	0.64
1:CA:7:G:H21	5:CE:121:LYS:CE	2.11	0.64
48:BZ:5:LYS:HB3	48:BZ:57:GLU:HB2	1.80	0.64
27:DE:102:PRO:HB2	27:DE:105:VAL:HG23	1.80	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
17:CQ:8:GLY:HA3	17:CQ:23:VAL:HG12	1.78	0.64
23:DA:663:G:H5''	34:DL:21:ARG:HD3	1.80	0.64
23:BA:1607:C:H4'	23:BA:1608:A:O5'	1.97	0.64
1:CA:323:U:H4'	20:CT:22:ARG:HB3	1.78	0.64
12:AL:26:LEU:CD1	12:AL:27:LYS:H	2.07	0.64
23:DA:81:G:H21	43:DU:2:ARG:NH2	1.96	0.64
23:BA:2873:A:N3	36:BN:6:SER:HB2	2.13	0.64
1:CA:986:A:H1'	19:CS:54:GLY:O	1.97	0.64
1:AA:1151:A:O2'	1:AA:1152:A:H8	1.79	0.64
11:CK:24:SER:HB3	11:CK:27:ASN:O	1.97	0.64
47:DY:6:VAL:O	47:DY:10:LEU:HG	1.97	0.64
1:AA:105:G:H2'	1:AA:106:C:H6	1.63	0.64
36:BN:9:LYS:O	36:BN:10:LEU:HG	1.96	0.64
12:AL:44:PRO:HG3	12:AL:52:ARG:HD3	1.79	0.64
23:DA:2271:G:OP1	45:DW:18:ALA:HB1	1.96	0.64
23:DA:1614:A:N6	41:DS:87:PRO:HA	2.11	0.64
10:CJ:55:LYS:HD2	10:CJ:55:LYS:O	1.96	0.64
42:BT:70:LEU:HD23	42:BT:71:GLY:N	2.13	0.64
35:BM:83:MET:O	35:BM:83:MET:HG3	1.98	0.64
45:BW:37:LEU:O	45:BW:38:VAL:HG23	1.98	0.64
32:DJ:135:LEU:HD23	32:DJ:136:GLY:N	2.12	0.64
32:BJ:36:TRP:HB2	32:BJ:156:GLN:CB	2.27	0.64
3:AC:22:TRP:HB3	3:AC:59:ARG:H	1.63	0.64
30:DH:82:ARG:HB3	30:DH:89:TYR:CB	2.27	0.64
24:DB:13:A:H5'	45:DW:74:ARG:HH21	1.62	0.64
23:DA:860:U:C5	23:DA:917:A:N7	2.66	0.64
1:AA:243:A:H4'	1:AA:244:U:O5'	1.97	0.64
1:AA:1194:U:H2'	1:AA:1195:C:C6	2.33	0.64
23:DA:2506:U:H5	23:DA:2583:G:H1	1.44	0.64
27:BE:31:HIS:HB2	34:BL:13:ASN:HB3	1.79	0.64
37:BO:33:LYS:O	37:BO:33:LYS:HD3	1.98	0.64
35:BM:110:THR:HB	35:BM:112:GLU:OE1	1.98	0.64
53:B5:39:LYS:O	53:B5:43:GLN:HG2	1.97	0.64
12:CL:44:PRO:HG3	12:CL:52:ARG:HD3	1.80	0.64
32:DJ:36:TRP:HB2	32:DJ:156:GLN:CB	2.28	0.64
29:BG:149:ARG:HA	29:BG:162:ILE:CG1	2.28	0.64
1:AA:1224:G:H4'	13:AM:102:ARG:HH22	1.62	0.64
29:DG:149:ARG:HA	29:DG:162:ILE:CG1	2.28	0.64
23:BA:1210:A:H4'	23:BA:1211:U:O5'	1.98	0.64
17:CQ:45:HIS:CD2	17:CQ:47:PRO:HD3	2.33	0.64
27:DE:139:PHE:CB	27:DE:166:ALA:HB1	2.28	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:AD:30:LYS:C	4:AD:32:ALA:H	2.00	0.64
23:BA:2599:G:C8	25:BC:237:GLU:HG3	2.32	0.64
1:CA:413:G:N2	1:CA:428:G:H1'	2.13	0.64
36:DN:38:VAL:HB	36:DN:39:PRO:HD3	1.78	0.64
36:DN:79:LEU:HD23	36:DN:83:ILE:HB	1.79	0.64
48:DZ:5:LYS:HB3	48:DZ:57:GLU:HB2	1.79	0.64
24:BB:79:C:H2'	24:BB:80:U:O4'	1.97	0.64
12:CL:23:VAL:HG13	12:CL:97:TYR:CE2	2.33	0.64
23:DA:2402:C:H5'	23:DA:2403:C:OP2	1.97	0.64
26:DD:51:PHE:H	26:DD:75:VAL:HB	1.60	0.64
36:DN:9:LYS:O	36:DN:10:LEU:HG	1.97	0.64
17:AQ:45:HIS:CD2	17:AQ:47:PRO:HD3	2.32	0.64
37:BO:31:SER:HB3	37:BO:34:HIS:HB2	1.78	0.64
23:BA:1529:A:H62	23:BA:1542:G:N2	1.96	0.64
25:DC:78:LYS:HD3	25:DC:114:GLY:HA2	1.80	0.64
1:CA:438:G:H4'	1:CA:439:A:OP1	1.97	0.64
1:CA:255:G:O3'	17:CQ:17:LYS:HD3	1.98	0.64
9:AI:28:VAL:HG22	9:AI:63:ILE:HB	1.79	0.64
23:BA:2777:G:H5''	23:BA:2778:A:H5'	1.78	0.64
36:BN:38:VAL:HB	36:BN:39:PRO:HD3	1.80	0.64
15:AO:48:LYS:HE2	15:AO:48:LYS:HA	1.78	0.64
23:BA:1007:C:O2'	32:BJ:131:PRO:HA	1.98	0.64
2:AB:168:THR:OG1	2:AB:192:SER:HA	1.98	0.64
1:AA:413:G:N2	1:AA:428:G:H1'	2.13	0.64
23:BA:664:C:H4'	23:BA:941:A:OP1	1.98	0.64
29:DG:101:ARG:NE	29:DG:101:ARG:H	1.93	0.64
25:BC:270:ILE:O	25:BC:271:ILE:HG13	1.97	0.64
1:CA:1298:C:N4	7:CG:114:ARG:HD3	2.13	0.64
9:AI:27:THR:O	9:AI:62:TYR:HA	1.98	0.64
23:BA:811:U:O2	23:BA:1250:G:H2'	1.98	0.64
23:BA:185:U:H2'	23:BA:186:G:C8	2.32	0.64
23:BA:480:A:OP2	43:BU:46:LYS:HE2	1.97	0.64
9:CI:28:VAL:HG22	9:CI:63:ILE:HB	1.79	0.64
43:BU:50:ARG:HD3	43:BU:51:VAL:H	1.63	0.64
25:BC:206:LEU:O	25:BC:211:ARG:HD3	1.98	0.64
23:BA:954:G:H5''	35:BM:13:GLN:CG	2.27	0.63
37:DO:66:ALA:HB1	37:DO:101:LEU:HD22	1.80	0.63
12:AL:70:PRO:HD3	12:AL:99:ILE:HG22	1.80	0.63
3:CC:59:ARG:HG2	3:CC:64:VAL:HG22	1.79	0.63
1:CA:265:G:H5'	17:CQ:64:PRO:O	1.97	0.63
32:DJ:112:LYS:O	32:DJ:116:THR:HG22	1.98	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:BN:79:LEU:HD23	36:BN:83:ILE:HB	1.79	0.63
23:BA:1434:A:H2'	23:BA:1435:G:C8	2.33	0.63
8:CH:64:LYS:HG2	8:CH:79:VAL:HG21	1.80	0.63
20:AT:26:ASN:HB2	20:AT:71:THR:HG23	1.79	0.63
23:DA:2687:U:C4	23:DA:2688:U:C5	2.86	0.63
23:BA:2271:G:OP1	45:BW:18:ALA:HB1	1.97	0.63
39:BQ:90:VAL:HG23	40:BR:39:LEU:HB3	1.81	0.63
23:BA:948:G:C8	23:BA:948:G:H5'	2.32	0.63
23:BA:2210:G:H21	23:BA:2211:G:H5'	1.64	0.63
12:CL:70:PRO:HD3	12:CL:99:ILE:HG22	1.81	0.63
37:DO:31:SER:HB3	37:DO:34:HIS:HB2	1.79	0.63
22:CV:6157:A:H61	22:CV:6172:U:H3	1.43	0.63
44:BV:24:LEU:HD21	44:BV:86:VAL:CG2	2.29	0.63
9:CI:27:THR:O	9:CI:62:TYR:HA	1.97	0.63
8:AH:91:ARG:HB2	12:AL:6:ILE:HD13	1.80	0.63
24:BB:11:C:H3'	24:BB:12:C:H6	1.63	0.63
26:DD:119:ARG:HD3	26:DD:120:TRP:CE2	2.33	0.63
23:DA:1803:A:H5''	23:DA:1804:C:OP2	1.98	0.63
1:AA:1126:U:H2'	1:AA:1127:G:C8	2.34	0.63
46:DX:86:SER:O	46:DX:90:ILE:HG12	1.98	0.63
23:DA:1694:C:C5'	23:DA:1694:C:H6	2.11	0.63
35:DM:83:MET:HG3	35:DM:83:MET:O	1.97	0.63
29:BG:101:ARG:N	29:BG:101:ARG:HE	1.94	0.63
46:BX:13:ILE:HD11	46:BX:15:ALA:HB2	1.79	0.63
46:BX:46:LEU:HD23	46:BX:46:LEU:O	1.99	0.63
23:DA:1657:C:H2'	23:DA:1658:C:C6	2.33	0.63
23:DA:126:A:OP2	52:D4:19:ARG:HB2	1.97	0.63
23:BA:1652:A:OP1	36:BN:9:LYS:HE3	1.98	0.63
23:BA:389:G:O6	34:BL:71:VAL:HG23	1.97	0.63
35:DM:6:ARG:O	35:DM:7:MET:HB2	1.98	0.63
23:BA:2294:C:H2'	23:BA:2295:C:H6	1.62	0.63
23:DA:1434:A:H2'	23:DA:1435:G:C8	2.34	0.63
4:CD:28:SER:HB3	4:CD:29:PRO:HD2	1.79	0.63
33:BK:103:ALA:HB1	33:BK:105:GLU:OE1	1.98	0.63
23:DA:140:A:H8	23:DA:1408:C:O2'	1.81	0.63
23:DA:586:A:H5'	27:DE:89:VAL:HG21	1.78	0.63
26:BD:119:ARG:HD3	26:BD:120:TRP:CE2	2.34	0.63
35:DM:20:ALA:HB1	35:DM:99:PRO:O	1.98	0.63
25:DC:158:ALA:HB3	25:DC:161:THR:HG21	1.80	0.63
2:CB:8:LYS:HA	2:CB:217:ARG:HH12	1.63	0.63
23:DA:637:A:O5'	34:DL:116:GLY:HA2	1.98	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BA:637:A:O5'	34:BL:116:GLY:HA2	1.99	0.63
3:CC:22:TRP:HB3	3:CC:59:ARG:H	1.63	0.63
25:DC:186:HIS:CD2	25:DC:188:GLU:H	2.16	0.63
23:DA:1937:A:O2'	23:DA:1938:A:H5'	1.98	0.63
6:CF:76:ALA:O	6:CF:80:ARG:HG2	1.98	0.63
39:DQ:92:ARG:HD3	39:DQ:94:ASN:HB3	1.79	0.63
1:CA:79:G:H2'	1:CA:80:G:C8	2.33	0.63
19:CS:49:ILE:H	19:CS:49:ILE:HD12	1.64	0.63
23:DA:2401:U:H2'	23:DA:2402:C:H5''	1.80	0.63
35:DM:110:THR:HB	35:DM:112:GLU:OE1	1.98	0.63
1:AA:1298:C:H4'	1:AA:1299:A:C8	2.33	0.63
23:BA:2401:U:H2'	23:BA:2402:C:H5''	1.79	0.63
8:CH:91:ARG:HB2	12:CL:6:ILE:HD13	1.80	0.63
21:AU:10:ARG:HA	21:AU:13:ILE:HB	1.80	0.63
2:CB:70:PHE:O	2:CB:92:TYR:HA	1.98	0.63
46:DX:19:GLN:HG2	46:DX:41:ARG:HB2	1.80	0.63
23:DA:2062:A:O2'	23:DA:2063:C:H5'	1.98	0.63
37:DO:24:LEU:HD13	37:DO:82:ILE:HG23	1.81	0.63
40:DR:39:LEU:HD12	40:DR:47:VAL:HG11	1.81	0.63
40:BR:39:LEU:HD12	40:BR:47:VAL:HG11	1.81	0.63
1:CA:674:G:H2'	1:CA:675:A:H8	1.64	0.63
23:DA:2873:A:N3	36:DN:6:SER:HB2	2.14	0.63
1:CA:38:G:N2	1:CA:397:A:H5'	2.13	0.63
23:DA:1210:A:H4'	23:DA:1211:U:O5'	1.99	0.63
23:BA:1314:C:C6	23:BA:1314:C:H5'	2.34	0.63
23:BA:1316:U:H2'	23:BA:1317:A:H8	1.63	0.63
1:AA:406:G:H5''	4:AD:5:ILE:HD12	1.81	0.63
9:CI:17:VAL:HA	9:CI:63:ILE:HG13	1.80	0.63
23:DA:2350:C:H5''	53:D5:42:ARG:HD3	1.81	0.63
23:DA:1437:C:H2'	23:DA:1438:U:C6	2.34	0.63
46:BX:86:SER:O	46:BX:90:ILE:HG12	1.97	0.63
25:DC:5:LYS:HD2	25:DC:5:LYS:N	2.14	0.63
37:DO:33:LYS:O	37:DO:33:LYS:HD3	1.97	0.63
16:AP:19:ILE:HG22	16:AP:36:ILE:HD11	1.80	0.63
27:BE:103:LYS:HA	27:BE:106:ARG:HG3	1.78	0.63
23:DA:7:G:H2'	23:DA:8:A:H8	1.64	0.63
25:BC:131:LEU:CD1	25:BC:136:ILE:HG12	2.29	0.63
11:AK:29:ILE:HG22	11:AK:44:SER:HB3	1.81	0.63
36:BN:18:LEU:HD11	36:BN:22:ARG:CZ	2.29	0.63
19:AS:29:ARG:HB2	19:AS:48:THR:H	1.64	0.63
23:DA:1694:C:H6	23:DA:1694:C:H5''	1.62	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:BL:14:LYS:O	34:BL:15:ARG:HB2	1.98	0.63
9:AI:113:LYS:H	9:AI:119:ALA:HA	1.61	0.63
26:DD:36:ARG:NH1	26:DD:86:PRO:HD2	2.14	0.63
2:AB:70:PHE:O	2:AB:92:TYR:HA	1.98	0.63
26:BD:36:ARG:NH1	26:BD:86:PRO:HD2	2.14	0.63
40:DR:38:LEU:O	40:DR:39:LEU:HD13	1.99	0.63
40:DR:40:LEU:HD23	40:DR:47:VAL:HG23	1.80	0.63
1:CA:1227:A:OP2	13:CM:111:LYS:HE3	1.99	0.63
46:DX:13:ILE:HD11	46:DX:15:ALA:HB2	1.80	0.63
1:AA:674:G:H2'	1:AA:675:A:H8	1.63	0.63
25:DC:61:LEU:O	25:DC:63:ARG:NH1	2.31	0.63
36:DN:10:LEU:CB	36:DN:17:ARG:HE	2.12	0.63
18:CR:70:ILE:O	18:CR:74:ARG:HG3	1.97	0.63
32:BJ:112:LYS:O	32:BJ:116:THR:HG22	1.99	0.63
23:BA:1694:C:C5'	23:BA:1694:C:H6	2.12	0.63
9:AI:69:GLY:O	9:AI:73:GLN:HG3	1.99	0.63
7:AG:126:ASP:HB3	7:AG:131:LYS:O	1.99	0.63
23:DA:2724:C:OP1	26:DD:118:LYS:HE3	1.98	0.63
23:BA:663:G:H5''	34:BL:21:ARG:HD3	1.79	0.63
23:BA:214:G:H1'	23:BA:216:A:O2'	1.99	0.63
3:AC:107:GLN:CD	3:AC:107:GLN:H	2.02	0.63
23:DA:1309:G:H3'	52:D4:9:ARG:HH12	1.64	0.63
29:DG:101:ARG:N	29:DG:101:ARG:HE	1.93	0.62
23:BA:780:G:H21	23:BA:783:A:N6	1.93	0.62
33:DK:71:ARG:NH1	38:DP:74:ARG:HH22	1.95	0.62
36:BN:17:ARG:HG3	36:BN:18:LEU:N	2.14	0.62
1:CA:392:G:H2'	1:CA:393:A:C8	2.32	0.62
25:DC:142:VAL:HG23	25:DC:192:THR:O	1.98	0.62
23:BA:2402:C:H5'	23:BA:2403:C:OP2	1.98	0.62
38:BP:50:ILE:HA	38:BP:99:LEU:HD11	1.81	0.62
1:CA:1071:C:H5''	5:CE:49:PRO:HG2	1.81	0.62
23:BA:1389:G:H2'	23:BA:1390:U:C6	2.34	0.62
2:AB:8:LYS:HA	2:AB:217:ARG:HH12	1.63	0.62
19:CS:11:VAL:HG23	19:CS:38:SER:HB2	1.81	0.62
23:DA:2777:G:H5''	23:DA:2778:A:H5'	1.80	0.62
23:DA:1388:G:H2'	23:DA:1389:G:H8	1.64	0.62
27:DE:170:LEU:HD12	27:DE:171:PRO:HD2	1.81	0.62
1:AA:976:G:H22	1:AA:136(A):C:H2'	1.64	0.62
43:DU:50:ARG:HD3	43:DU:51:VAL:H	1.64	0.62
34:BL:62:LEU:N	34:BL:62:LEU:HD13	2.14	0.62
37:BO:66:ALA:HB1	37:BO:101:LEU:HD22	1.80	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
52:D4:8:ASN:C	52:D4:8:ASN:ND2	2.53	0.62
23:DA:1314:C:C6	23:DA:1314:C:H5'	2.34	0.62
19:AS:18:LYS:O	19:AS:22:LEU:HD23	1.99	0.62
19:AS:49:ILE:HD12	19:AS:49:ILE:H	1.63	0.62
23:DA:65:C:H2'	23:DA:66:C:C6	2.34	0.62
1:CA:1298:C:H41	7:CG:114:ARG:HD3	1.64	0.62
29:BG:84:SER:HA	29:BG:133:VAL:O	1.99	0.62
5:CE:82:VAL:HG21	5:CE:138:ALA:HA	1.80	0.62
23:BA:1953:A:H2	23:BA:2549:G:N3	1.97	0.62
25:DC:95:LEU:O	25:DC:95:LEU:HD12	1.98	0.62
23:BA:510:C:H2'	23:BA:511:U:O4'	1.99	0.62
1:CA:137:C:O4'	16:CP:63:GLY:HA3	1.99	0.62
23:BA:242:G:C8	53:B5:5:LYS:HG2	2.33	0.62
27:BE:29:ASN:H	27:BE:112:MET:HE1	1.64	0.62
23:DA:747:U:OP2	50:D2:3:LYS:HD3	1.99	0.62
7:AG:111:ARG:HB3	7:AG:113:GLU:HG2	1.80	0.62
1:CA:979:C:H42	14:CN:18:VAL:HG12	1.63	0.62
1:AA:427:U:OP1	4:AD:13:ARG:NH2	2.32	0.62
11:AK:24:SER:HB3	11:AK:27:ASN:O	1.99	0.62
25:BC:10:THR:HG23	25:BC:13:ARG:CB	2.29	0.62
27:DE:31:HIS:HB2	34:DL:13:ASN:HB3	1.80	0.62
34:DL:14:LYS:O	34:DL:15:ARG:HB2	1.99	0.62
21:CU:10:ARG:HA	21:CU:13:ILE:HB	1.80	0.62
3:CC:107:GLN:H	3:CC:107:GLN:CD	2.03	0.62
6:CF:12:PRO:HD3	6:CF:58:GLY:HA2	1.81	0.62
32:BJ:135:LEU:HD23	32:BJ:136:GLY:N	2.13	0.62
33:BK:71:ARG:NH1	38:BP:74:ARG:HH22	1.98	0.62
34:DL:30:THR:HG22	34:DL:31:ALA:N	2.15	0.62
23:DA:1316:U:H2'	23:DA:1317:A:H8	1.64	0.62
23:DA:1389:G:H2'	23:DA:1390:U:C6	2.34	0.62
23:DA:2078:C:H2'	23:DA:2079:U:H6	1.64	0.62
34:BL:9:ASN:N	34:BL:10:PRO:HD3	2.15	0.62
23:BA:2687:U:C4	23:BA:2688:U:C5	2.88	0.62
42:DT:64:LYS:HG2	42:DT:65:ARG:H	1.64	0.62
30:BH:6:LEU:HD23	30:BH:6:LEU:H	1.63	0.62
23:DA:1022:G:H8	32:DJ:92:GLN:NE2	1.96	0.62
34:DL:62:LEU:HD13	34:DL:62:LEU:N	2.14	0.62
26:BD:132:HIS:CD2	26:BD:135:HIS:NE2	2.68	0.62
2:CB:27:LYS:HG3	2:CB:194:PRO:HD2	1.82	0.62
23:BA:2075:U:H2'	23:BA:2238:G:H22	1.63	0.62
9:AI:17:VAL:HA	9:AI:63:ILE:HG13	1.80	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:DC:206:LEU:O	25:DC:211:ARG:HD3	2.00	0.62
25:DC:79:VAL:HG21	25:DC:111:LEU:HD11	1.80	0.62
1:CA:1349:A:H2'	1:CA:1350:A:C8	2.35	0.62
27:DE:178:PRO:HB3	27:DE:201:VAL:HG11	1.82	0.62
23:BA:2502:G:H5'	23:BA:2503:A:H5''	1.81	0.62
33:DK:20:MET:HG2	33:DK:21:CYS:O	2.00	0.62
33:BK:104:ARG:HH11	33:BK:104:ARG:HB3	1.63	0.62
6:AF:60:PHE:C	6:AF:61:LEU:HD12	2.19	0.62
39:BQ:21:ALA:CB	39:BQ:35:ALA:HB1	2.29	0.62
23:BA:2350:C:H5''	53:B5:42:ARG:HD3	1.81	0.62
7:CG:111:ARG:HB3	7:CG:113:GLU:HG2	1.80	0.62
39:DQ:90:VAL:HG23	40:DR:39:LEU:HB3	1.81	0.62
1:CA:922:G:H2'	1:CA:923:A:H8	1.63	0.62
42:DT:63:LYS:HZ1	42:DT:72:LYS:HB3	1.63	0.62
26:DD:9:VAL:HG22	26:DD:25:VAL:HB	1.81	0.62
34:BL:16:ARG:NH1	34:BL:18:ARG:HG3	2.15	0.62
23:DA:1794:U:H2'	23:DA:1795:C:C6	2.35	0.62
44:BV:102:LEU:HD23	44:BV:137:ILE:HB	1.81	0.62
23:BA:185:U:H4'	23:BA:218:A:H4'	1.82	0.62
1:AA:976:G:N2	1:AA:136(A):C:H2'	2.14	0.62
24:BB:50:G:OP2	37:BO:62:LYS:HD3	2.00	0.62
46:DX:25:LYS:HG2	46:DX:35:THR:HG22	1.80	0.62
32:DJ:127:LYS:HB2	32:DJ:140:PHE:CE1	2.33	0.62
4:AD:166:LYS:HE2	25:DC:134:ARG:HH21	1.63	0.62
2:CB:178:ARG:HE	8:CH:74:PRO:HD3	1.65	0.62
23:DA:2210:G:H21	23:DA:2211:G:H5'	1.64	0.62
19:CS:18:LYS:O	19:CS:22:LEU:HD23	2.00	0.62
1:AA:190:G:H4'	1:AA:191(A):G:OP2	2.00	0.62
9:AI:26:VAL:HG13	9:AI:61:ALA:HB3	1.82	0.62
43:DU:90:LEU:HG	43:DU:91:GLU:N	2.15	0.62
1:AA:948:C:OP1	13:AM:107:ALA:HA	1.99	0.62
7:CG:126:ASP:HB3	7:CG:131:LYS:O	2.00	0.62
23:DA:2307:G:H2'	23:DA:2308:G:H5'	1.81	0.62
33:DK:104:ARG:HH11	33:DK:104:ARG:HB3	1.65	0.62
45:DW:23:VAL:HA	45:DW:38:VAL:CG2	2.24	0.62
23:DA:1529:A:H62	23:DA:1542:G:N2	1.98	0.62
33:DK:2:ILE:HD11	33:DK:82:ASN:HD22	1.64	0.62
44:DV:69:THR:HG22	44:DV:90:VAL:HG22	1.82	0.62
29:DG:68:THR:O	29:DG:72:ILE:HG12	2.00	0.62
23:BA:492:A:H2'	23:BA:493:G:O4'	2.00	0.62
46:DX:46:LEU:HD23	46:DX:46:LEU:O	2.00	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:624:C:O3'	16:AP:10:GLY:HA2	2.00	0.62
35:BM:38:GLU:HB2	35:BM:127:ILE:HG23	1.81	0.62
45:BW:23:VAL:HA	45:BW:38:VAL:CG2	2.24	0.62
43:BU:7:VAL:HB	43:BU:8:LYS:HZ2	1.65	0.62
23:BA:81:G:H21	43:BU:2:ARG:NH2	1.98	0.62
23:BA:343:C:C6	23:BA:343:C:H5'	2.32	0.62
17:CQ:64:PRO:HA	17:CQ:70:ARG:HG3	1.82	0.62
30:BH:78:THR:HA	30:BH:143:SER:HB3	1.80	0.62
5:AE:6:PHE:CD2	5:AE:36:ASP:HB3	2.35	0.62
23:DA:1542:G:H4'	23:DA:1543:A:O5'	2.00	0.62
23:BA:674:G:H2'	23:BA:804:A:H61	1.62	0.62
4:AD:22:LYS:HB2	4:AD:26:CYS:SG	2.40	0.62
44:DV:102:LEU:HD23	44:DV:137:ILE:HB	1.80	0.62
17:CQ:40:LYS:HD2	17:CQ:42:TYR:CZ	2.35	0.62
46:BX:73:LEU:HD21	46:BX:94:LEU:HG	1.82	0.62
42:DT:59:VAL:HB	42:DT:76:ARG:HG3	1.81	0.62
23:BA:2577:A:H5''	23:BA:2578:G:H5'	1.81	0.62
19:AS:62:ILE:HA	19:AS:66:MET:SD	2.40	0.62
29:DG:84:SER:HA	29:DG:133:VAL:O	2.00	0.62
1:CA:370:C:H2'	1:CA:371:G:H8	1.65	0.62
23:BA:1937:A:O2'	23:BA:1938:A:H5'	1.99	0.62
46:BX:25:LYS:HG2	46:BX:35:THR:HG22	1.81	0.62
23:BA:2186:G:H2'	23:BA:2187:G:H8	1.65	0.62
39:DQ:92:ARG:HD2	39:DQ:95:LEU:HG	1.81	0.62
26:DD:132:HIS:HA	26:DD:135:HIS:NE2	2.15	0.62
25:BC:30:GLU:HG3	25:BC:63:ARG:HH21	1.63	0.62
5:CE:6:PHE:CD2	5:CE:36:ASP:HB3	2.35	0.62
1:CA:1220:G:H21	19:CS:54:GLY:CA	2.13	0.62
34:BL:30:THR:HG22	34:BL:31:ALA:N	2.15	0.62
46:DX:27:GLU:HB3	46:DX:33:LYS:HG3	1.82	0.62
1:AA:946:A:H2'	1:AA:947:G:H8	1.65	0.62
23:BA:273(G):C:H2'	23:BA:274:G:H5''	1.82	0.62
39:DQ:21:ALA:CB	39:DQ:35:ALA:HB1	2.30	0.62
23:DA:38:A:H2'	23:DA:39:C:C6	2.34	0.62
32:BJ:127:LYS:HB2	32:BJ:140:PHE:CE1	2.35	0.62
23:DA:492:A:H2'	23:DA:493:G:O4'	2.00	0.62
6:CF:60:PHE:C	6:CF:61:LEU:HD12	2.20	0.62
23:DA:664:C:H4'	23:DA:941:A:OP1	2.00	0.61
23:BA:1046:A:H1'	31:BI:4:LYS:HD3	1.82	0.61
50:D2:40:LYS:CE	50:D2:46:CYS:HB3	2.29	0.61
19:CS:31:ILE:HG23	19:CS:49:ILE:HA	1.81	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DA:1343:G:H8	23:DA:1343:G:H5'	1.64	0.61
4:CD:30:LYS:C	4:CD:32:ALA:H	2.03	0.61
46:DX:27:GLU:CB	46:DX:33:LYS:HA	2.30	0.61
13:AM:14:ARG:CZ	13:AM:42:ALA:HA	2.30	0.61
23:DA:779:U:H5'	25:DC:49:ILE:HD11	1.82	0.61
23:DA:1007:C:O2'	32:DJ:131:PRO:HA	1.99	0.61
35:BM:20:ALA:HB1	35:BM:99:PRO:O	2.00	0.61
1:CA:801:U:H2'	1:CA:802:A:C8	2.35	0.61
1:CA:1126:U:H2'	1:CA:1127:G:C8	2.35	0.61
6:AF:12:PRO:HD3	6:AF:58:GLY:HA2	1.82	0.61
24:DB:50:G:OP2	37:DO:62:LYS:HD3	2.00	0.61
12:AL:23:VAL:HG13	12:AL:97:TYR:CE2	2.35	0.61
1:AA:801:U:H2'	1:AA:802:A:C8	2.35	0.61
1:CA:1322:C:OP2	13:CM:100:GLY:HA3	2.00	0.61
26:BD:4:ILE:CG1	26:BD:28:ALA:HB1	2.30	0.61
43:DU:8:LYS:HE2	43:DU:37:VAL:HG11	1.81	0.61
23:BA:1021:A:O2'	23:BA:1123:C:H5''	2.00	0.61
23:DA:2393:A:C5'	34:DL:62:LEU:HD12	2.31	0.61
43:BU:8:LYS:HE2	43:BU:37:VAL:HG11	1.82	0.61
30:BH:123:LEU:HD23	30:BH:124:GLY:N	2.15	0.61
2:AB:25:ASN:HD22	2:AB:25:ASN:N	1.98	0.61
23:DA:860:U:O2'	23:DA:861:A:H5'	1.99	0.61
1:AA:392:G:H2'	1:AA:393:A:C8	2.34	0.61
34:DL:71:VAL:HB	34:DL:72:PRO:HD3	1.82	0.61
46:DX:19:GLN:HE21	46:DX:41:ARG:HE	1.48	0.61
23:DA:2787:C:H1'	26:DD:62:PRO:HG3	1.81	0.61
23:BA:2653:U:H3	23:BA:2667:C:H42	1.48	0.61
23:BA:729:G:C5	25:BC:208:LYS:HB2	2.35	0.61
25:BC:21:PHE:O	25:BC:24:ILE:HG22	2.00	0.61
23:DA:2502:G:H5'	23:DA:2503:A:H5''	1.81	0.61
1:AA:754:C:O2	1:AA:754:C:H3'	2.01	0.61
27:BE:199:TRP:O	27:BE:203:GLN:HG2	2.00	0.61
23:DA:2346:A:H5''	23:DA:2383:G:H1'	1.82	0.61
32:DJ:57:LEU:HD21	32:DJ:143:LEU:HB2	1.82	0.61
30:DH:126:TYR:H	30:DH:142:VAL:HB	1.65	0.61
1:AA:370:C:H2'	1:AA:371:G:H8	1.66	0.61
23:BA:1291:C:H2'	23:BA:1292:U:C6	2.35	0.61
33:BK:2:ILE:HD11	33:BK:82:ASN:HD22	1.65	0.61
25:BC:108:PRO:HG3	25:BC:143:HIS:CE1	2.36	0.61
9:CI:79:LEU:HD23	9:CI:101:PHE:O	1.99	0.61
1:CA:600:C:OP1	8:CH:97:VAL:HG12	2.00	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:DD:132:HIS:CD2	26:DD:135:HIS:NE2	2.68	0.61
32:BJ:57:LEU:HD21	32:BJ:143:LEU:HB2	1.80	0.61
34:DL:16:ARG:NH1	34:DL:18:ARG:HG3	2.16	0.61
6:CF:37:VAL:HA	6:CF:65:VAL:HG12	1.81	0.61
16:CP:19:ILE:HG22	16:CP:36:ILE:HD11	1.82	0.61
23:BA:2562:U:H1'	33:BK:23:ARG:NH1	2.15	0.61
23:DA:729:G:C5	25:DC:208:LYS:HB2	2.36	0.61
33:DK:47:ILE:HG13	33:DK:48:PRO:HD2	1.82	0.61
9:CI:26:VAL:HG13	9:CI:61:ALA:HB3	1.81	0.61
24:DB:11:C:H3'	24:DB:12:C:H6	1.65	0.61
23:DA:639:U:H2'	23:DA:640:C:C6	2.35	0.61
42:BT:64:LYS:HG2	42:BT:65:ARG:H	1.65	0.61
23:DA:2593:U:H2'	23:DA:2594:C:C6	2.36	0.61
35:DM:38:GLU:O	35:DM:127:ILE:HD13	2.01	0.61
35:DM:38:GLU:HB2	35:DM:127:ILE:HG23	1.83	0.61
17:AQ:40:LYS:HD2	17:AQ:42:TYR:CZ	2.35	0.61
10:AJ:48:THR:HA	10:AJ:62:HIS:CB	2.29	0.61
23:DA:1408:C:C2	23:DA:1595:G:N2	2.69	0.61
36:DN:17:ARG:HG3	36:DN:18:LEU:N	2.14	0.61
48:DZ:1:MET:SD	48:DZ:40:THR:HA	2.41	0.61
4:CD:4:TYR:HE1	4:CD:11:LEU:HD11	1.65	0.61
4:AD:166:LYS:HE2	25:DC:134:ARG:NH2	2.15	0.61
38:BP:98:LYS:HB3	38:BP:100:TYR:CE1	2.36	0.61
43:DU:76:CYS:HB3	43:DU:77:PRO:HD2	1.81	0.61
23:BA:639:U:H2'	23:BA:640:C:C6	2.35	0.61
38:DP:41:ARG:HD2	38:DP:42:ILE:H	1.66	0.61
38:DP:50:ILE:HA	38:DP:99:LEU:HD11	1.82	0.61
23:BA:7:G:H2'	23:BA:8:A:H8	1.64	0.61
23:DA:2294:C:H2'	23:DA:2295:C:H6	1.66	0.61
1:AA:684:A:H2'	1:AA:685:G:C8	2.35	0.61
11:AK:120:ARG:HH21	11:AK:126:ARG:HH21	1.49	0.61
25:BC:5:LYS:N	25:BC:5:LYS:HD2	2.14	0.61
27:BE:102:PRO:HB2	27:BE:105:VAL:HG23	1.81	0.61
1:AA:1349:A:H2'	1:AA:1350:A:C8	2.35	0.61
23:DA:774:A:H2	23:DA:787:U:HO2'	1.47	0.61
5:AE:82:VAL:HG21	5:AE:138:ALA:HA	1.81	0.61
23:BA:996:A:C4'	39:BQ:92:ARG:HH12	2.12	0.61
23:DA:780:G:H21	23:DA:783:A:N6	1.94	0.61
47:DY:2:LYS:HD2	47:DY:2:LYS:H	1.65	0.61
50:D2:40:LYS:HE2	50:D2:46:CYS:HB3	1.81	0.61
25:DC:44:ASN:HB3	25:DC:50:THR:HG21	1.82	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:BC:25:THR:HG22	25:BC:82:ILE:O	2.01	0.61
4:AD:4:TYR:HE1	4:AD:11:LEU:HD11	1.65	0.61
23:BA:1678:G:N2	23:BA:1989:G:H22	1.98	0.61
23:BA:860:U:C5	23:BA:917:A:N7	2.68	0.61
1:AA:920:U:H2'	1:AA:921:U:C6	2.36	0.61
35:BM:6:ARG:O	35:BM:7:MET:HB2	2.00	0.61
8:AH:64:LYS:HG2	8:AH:79:VAL:HG21	1.82	0.61
41:DS:78:GLU:OE2	41:DS:99:ARG:HD3	2.00	0.61
27:BE:83:PHE:O	27:BE:84:VAL:C	2.38	0.61
29:BG:55:PRO:HG2	29:BG:61:HIS:HD2	1.63	0.61
1:AA:932:C:OP1	7:AG:4:ARG:HG2	2.01	0.61
1:AA:114:U:H2'	1:AA:115:G:C8	2.36	0.61
36:DN:84:ALA:HB3	36:DN:85:PRO:HD3	1.83	0.61
42:DT:70:LEU:HD23	42:DT:71:GLY:N	2.13	0.61
18:AR:26:LEU:HD13	18:AR:39:VAL:HG13	1.83	0.61
38:BP:41:ARG:HD2	38:BP:42:ILE:H	1.65	0.61
1:CA:737:A:H2'	1:CA:738:C:C6	2.35	0.61
11:CK:29:ILE:HG22	11:CK:44:SER:HB3	1.82	0.61
23:BA:245:G:H2'	23:BA:246:C:H6	1.65	0.61
9:CI:69:GLY:O	9:CI:73:GLN:HG3	2.00	0.61
36:DN:4:LEU:HG	36:DN:4:LEU:O	2.00	0.61
25:DC:267:SER:O	25:DC:270:ILE:HG13	2.01	0.61
34:DL:140:ALA:O	34:DL:141:ALA:HB2	2.00	0.61
23:BA:1408:C:C2	23:BA:1595:G:N2	2.68	0.61
28:DF:10:LYS:O	28:DF:14:GLU:HB3	2.00	0.61
23:BA:1388:G:H2'	23:BA:1389:G:H8	1.64	0.61
5:CE:101:ILE:HD11	5:CE:119:LEU:HD23	1.82	0.61
46:BX:67:ILE:N	46:BX:68:PRO:HD2	2.15	0.61
23:BA:2850:A:H5'	23:BA:2850:A:C8	2.36	0.61
23:DA:2562:U:H1'	33:DK:23:ARG:NH1	2.15	0.61
38:DP:98:LYS:HB3	38:DP:100:TYR:CE1	2.36	0.61
26:DD:4:ILE:CG1	26:DD:28:ALA:HB1	2.31	0.61
41:BS:78:GLU:OE2	41:BS:99:ARG:HD3	2.00	0.61
45:BW:21:LEU:H	45:BW:21:LEU:HD12	1.66	0.61
23:BA:2787:C:H1'	26:BD:62:PRO:HG3	1.82	0.61
23:DA:2393:A:H5''	34:DL:62:LEU:HB3	1.83	0.61
1:CA:1347:G:N2	1:CA:1373:G:H2'	2.14	0.61
25:BC:31:LYS:O	25:BC:36:PRO:HD3	2.01	0.61
25:DC:30:GLU:HG3	25:DC:63:ARG:HH21	1.64	0.61
34:BL:140:ALA:O	34:BL:141:ALA:HB2	2.00	0.61
23:DA:1678:G:N2	23:DA:1989:G:H22	1.98	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:BD:51:PHE:CD1	26:BD:52:LEU:HG	2.36	0.61
23:DA:1543:A:H5'	23:DA:1544:C:OP2	2.00	0.61
1:AA:1347:G:N2	1:AA:1373:G:H2'	2.15	0.61
23:BA:1358:G:O2'	23:BA:1359:A:H5''	2.01	0.61
26:BD:32:PRO:HA	26:BD:90:THR:HG22	1.83	0.61
42:BT:59:VAL:HB	42:BT:76:ARG:HG3	1.81	0.61
34:DL:91:PHE:CE2	34:DL:95:VAL:HG12	2.35	0.61
26:BD:173:VAL:HG12	26:BD:174:ASP:H	1.65	0.61
32:BJ:62:ARG:NH2	32:BJ:64:ASP:HB2	2.15	0.61
1:CA:475:G:H2'	1:CA:476:G:H8	1.66	0.61
25:DC:21:PHE:O	25:DC:24:ILE:HG22	2.01	0.61
32:DJ:62:ARG:NH2	32:DJ:64:ASP:HB2	2.16	0.61
43:BU:90:LEU:HG	43:BU:91:GLU:N	2.16	0.61
23:DA:1021:A:O2'	23:DA:1123:C:H5''	2.01	0.61
25:DC:25:THR:HG22	25:DC:82:ILE:O	2.00	0.61
12:AL:70:PRO:HD2	12:AL:101:ARG:HD3	1.82	0.61
2:AB:27:LYS:HG3	2:AB:194:PRO:HD2	1.81	0.61
23:DA:674:G:H2'	23:DA:804:A:H61	1.66	0.61
23:DA:1429:G:H2'	23:DA:1430:C:H6	1.65	0.61
19:AS:31:ILE:HG23	19:AS:49:ILE:HA	1.81	0.61
23:BA:1542:G:H4'	23:BA:1543:A:O5'	2.01	0.61
1:CA:1152:A:OP1	10:CJ:68:HIS:CD2	2.54	0.61
43:DU:76:CYS:CB	43:DU:77:PRO:HD2	2.31	0.61
23:BA:38:A:H2'	23:BA:39:C:C6	2.35	0.61
25:DC:70:TRP:CZ2	25:DC:150:LYS:HA	2.36	0.61
41:BS:65:LEU:HB2	41:BS:68:ARG:HG2	1.82	0.61
11:CK:57:THR:HG22	11:CK:59:TYR:H	1.64	0.61
41:DS:65:LEU:HB2	41:DS:68:ARG:HG2	1.83	0.61
47:DY:17:SER:HB3	47:DY:18:PRO:CD	2.30	0.61
23:BA:245:G:O6	53:B5:8:LYS:HE3	2.01	0.61
23:BA:1311:G:C5'	23:BA:1311:G:H8	2.12	0.61
30:DH:83:ALA:HB3	30:DH:123:LEU:HD12	1.82	0.61
30:BH:126:TYR:H	30:BH:142:VAL:HB	1.65	0.61
23:BA:65:C:H2'	23:BA:66:C:C6	2.35	0.61
6:AF:37:VAL:HA	6:AF:65:VAL:HG12	1.82	0.61
44:DV:58:VAL:HA	44:DV:67:LEU:O	2.01	0.61
41:DS:29:LEU:HD22	41:DS:69:LEU:HD11	1.81	0.61
19:AS:11:VAL:HG23	19:AS:38:SER:HB2	1.81	0.61
53:B5:57:ARG:HB2	53:B5:57:ARG:HH11	1.66	0.60
27:BE:53:THR:HG22	27:BE:56:GLU:CD	2.22	0.60
23:BA:2039:C:H2'	23:BA:2040:C:C6	2.35	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DA:2075:U:H2'	23:DA:2238:G:H22	1.64	0.60
23:BA:1543:A:H5'	23:BA:1544:C:OP2	2.00	0.60
46:BX:27:GLU:CB	46:BX:33:LYS:HA	2.31	0.60
23:DA:185:U:H2'	23:DA:186:G:H8	1.63	0.60
25:BC:69:ARG:HH21	25:BC:130:ALA:HB2	1.62	0.60
23:BA:2078:C:H2'	23:BA:2079:U:C6	2.36	0.60
28:BF:36:LYS:HB3	28:BF:160:VAL:HB	1.83	0.60
23:DA:2186:G:H2'	23:DA:2187:G:H8	1.65	0.60
27:DE:83:PHE:O	27:DE:84:VAL:C	2.40	0.60
1:CA:1104:G:H4'	2:CB:111:ARG:NH1	2.15	0.60
1:CA:754:C:H3'	1:CA:754:C:O2	2.00	0.60
4:CD:63:LYS:HD2	4:CD:198:VAL:HG22	1.83	0.60
23:BA:1437:C:H2'	23:BA:1438:U:C6	2.36	0.60
1:AA:735:C:H2'	1:AA:736:C:H6	1.66	0.60
52:B4:8:ASN:C	52:B4:8:ASN:ND2	2.54	0.60
1:AA:1220:G:H21	19:AS:54:GLY:CA	2.14	0.60
46:BX:31:GLY:O	46:BX:32:LYS:HB2	2.01	0.60
23:BA:1478:G:O2'	23:BA:1558:A:H2	1.84	0.60
36:BN:84:ALA:HB3	36:BN:85:PRO:HD3	1.83	0.60
29:BG:68:THR:O	29:BG:72:ILE:HG12	2.01	0.60
23:DA:242:G:C8	53:D5:5:LYS:HG2	2.35	0.60
27:DE:199:TRP:O	27:DE:203:GLN:HG2	2.00	0.60
27:DE:164:ARG:HG3	27:DE:175:THR:OG1	2.00	0.60
4:AD:63:LYS:HD2	4:AD:198:VAL:HG22	1.83	0.60
3:AC:66:VAL:HB	3:AC:101:LEU:HD23	1.84	0.60
2:AB:87:ARG:O	2:AB:87:ARG:HD2	2.01	0.60
35:DM:30:GLY:HA2	35:DM:107:ALA:HB2	1.83	0.60
23:DA:1587:A:H2'	23:DA:1588:C:C6	2.37	0.60
23:BA:1309:G:H3'	52:B4:9:ARG:HH12	1.65	0.60
32:DJ:80:ALA:O	32:DJ:83:ILE:HG13	2.01	0.60
23:BA:747:U:OP2	50:B2:3:LYS:HD3	2.00	0.60
23:DA:996:A:C4'	39:DQ:92:ARG:HH12	2.13	0.60
14:AN:32:SER:HB3	14:AN:41:ARG:HG2	1.84	0.60
43:BU:76:CYS:HB3	43:BU:77:PRO:HD2	1.83	0.60
36:BN:4:LEU:O	36:BN:4:LEU:HG	2.01	0.60
53:B5:22:VAL:HB	53:B5:54:GLU:CG	2.31	0.60
1:AA:922:G:H2'	1:AA:923:A:H8	1.64	0.60
25:BC:61:LEU:O	25:BC:63:ARG:NH1	2.34	0.60
36:BN:10:LEU:CB	36:BN:17:ARG:HE	2.13	0.60
23:BA:2074:U:H2'	23:BA:2075:U:C6	2.36	0.60
19:CS:6:LYS:HD3	19:CS:7:LYS:HE2	1.83	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DA:795:C:H2'	23:DA:796:C:H6	1.65	0.60
23:DA:1953:A:H2	23:DA:2549:G:N3	1.98	0.60
1:AA:475:G:H2'	1:AA:476:G:H8	1.65	0.60
1:CA:33:A:H2'	1:CA:34:C:C6	2.36	0.60
23:DA:85:G:H5''	23:DA:85:G:H8	1.65	0.60
1:AA:7:G:H21	5:AE:121:LYS:HE2	1.67	0.60
30:DH:90:GLY:O	30:DH:91:SER:HB2	2.00	0.60
23:DA:214:G:H1'	23:DA:216:A:O2'	2.00	0.60
33:DK:3:GLN:CB	33:DK:4:PRO:HD2	2.27	0.60
53:D5:22:VAL:HB	53:D5:54:GLU:CG	2.31	0.60
46:BX:19:GLN:HG2	46:BX:41:ARG:HB2	1.83	0.60
23:DA:510:C:H2'	23:DA:511:U:O4'	2.00	0.60
10:CJ:24:VAL:HG21	10:CJ:37:PRO:HD3	1.82	0.60
13:AM:27:LYS:HG3	13:AM:31:LYS:HE3	1.83	0.60
23:DA:2850:A:C8	23:DA:2850:A:H5'	2.36	0.60
23:BA:2893:G:H5''	23:BA:2894:G:O4'	2.02	0.60
42:DT:8:ILE:HD12	42:DT:8:ILE:H	1.64	0.60
5:CE:47:LYS:N	5:CE:47:LYS:HD3	2.16	0.60
23:BA:85:G:H8	23:BA:85:G:H5''	1.66	0.60
7:AG:46:ALA:O	7:AG:50:ILE:HG12	2.02	0.60
1:CA:1528:U:H6	1:CA:1528:U:H5''	1.66	0.60
25:DC:34:VAL:O	25:DC:35:LYS:HD3	2.01	0.60
25:BC:34:VAL:O	25:BC:35:LYS:HD3	2.01	0.60
26:DD:51:PHE:CD1	26:DD:52:LEU:HG	2.36	0.60
46:BX:83:GLU:HG2	46:BX:84:GLY:N	2.17	0.60
42:BT:8:ILE:N	42:BT:8:ILE:HD12	2.16	0.60
46:DX:73:LEU:HD21	46:DX:94:LEU:HG	1.83	0.60
18:CR:26:LEU:HD13	18:CR:39:VAL:HG13	1.83	0.60
2:AB:178:ARG:HE	8:AH:74:PRO:HD3	1.66	0.60
23:DA:2653:U:H3	23:DA:2667:C:H42	1.50	0.60
1:AA:299:G:H2'	1:AA:300:A:C8	2.36	0.60
1:CA:920:U:H2'	1:CA:921:U:C6	2.36	0.60
5:AE:48:ALA:HB2	5:AE:57:LYS:HD3	1.84	0.60
2:CB:87:ARG:O	2:CB:87:ARG:HD2	2.02	0.60
1:CA:505:G:H2'	1:CA:506:G:H8	1.66	0.60
34:BL:38:GLN:HG3	34:BL:39:LYS:H	1.66	0.60
23:BA:2439:A:C5'	23:BA:2439:A:C8	2.84	0.60
1:AA:1346:A:H5'	9:AI:120:ARG:NH1	2.16	0.60
11:CK:29:ILE:HG22	11:CK:44:SER:CB	2.31	0.60
34:DL:57:THR:HG23	34:DL:59:LEU:CB	2.31	0.60
1:CA:1014:A:C2	1:CA:1219:U:H1'	2.37	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DA:342:G:H2'	23:DA:343:C:H5''	1.82	0.60
23:BA:558:G:OP1	32:BJ:134:PRO:HD2	2.00	0.60
28:BF:10:LYS:O	28:BF:14:GLU:HB3	2.00	0.60
23:BA:795:C:H2'	23:BA:796:C:H6	1.66	0.60
15:CO:5:LYS:H	15:CO:5:LYS:HD3	1.66	0.60
1:AA:1014:A:C2	1:AA:1219:U:H1'	2.37	0.60
32:BJ:117:HIS:CE1	32:BJ:120:ARG:HE	2.19	0.60
44:BV:24:LEU:HD21	44:BV:86:VAL:HG21	1.82	0.60
28:BF:114:ILE:HG23	28:BF:115:ARG:HD2	1.84	0.60
1:AA:1106:G:H5''	3:AC:172:ARG:HG2	1.83	0.60
36:BN:11:ASN:OD1	36:BN:12:ARG:N	2.30	0.60
9:AI:79:LEU:HD23	9:AI:101:PHE:O	2.00	0.60
26:DD:32:PRO:HA	26:DD:90:THR:HG22	1.83	0.60
23:BA:1771:C:HO2'	23:BA:1786:A:H8	1.50	0.60
10:AJ:24:VAL:HG21	10:AJ:37:PRO:HD3	1.83	0.60
23:DA:83:G:H22	23:DA:102:G:H2'	1.65	0.60
1:CA:1223:C:C5'	1:CA:1224:G:H5''	2.30	0.60
23:DA:955:C:OP1	35:DM:85:LYS:HE2	2.00	0.60
26:BD:132:HIS:HA	26:BD:135:HIS:NE2	2.15	0.60
23:DA:1568:G:P	25:DC:63:ARG:HH22	2.24	0.60
46:DX:31:GLY:O	46:DX:32:LYS:HB2	2.01	0.60
1:AA:1152:A:OP1	10:AJ:68:HIS:HD2	1.85	0.60
19:AS:5:LEU:HD12	19:AS:8:GLY:O	2.02	0.60
5:AE:101:ILE:HD11	5:AE:119:LEU:HD23	1.83	0.60
23:DA:1156:A:H4'	23:DA:1157:G:OP2	2.02	0.60
46:DX:67:ILE:N	46:DX:68:PRO:HD2	2.15	0.60
27:BE:164:ARG:HG3	27:BE:175:THR:OG1	2.02	0.60
19:CS:62:ILE:HA	19:CS:66:MET:SD	2.41	0.60
23:BA:153:C:OP1	46:BX:92:LYS:HE2	2.02	0.60
10:CJ:6:ILE:HG12	10:CJ:72:VAL:O	2.01	0.60
7:CG:26:PHE:O	7:CG:30:ILE:HG12	2.01	0.60
1:CA:328:C:H4'	1:CA:329:A:C5'	2.31	0.60
43:DU:71:LYS:HZ2	43:DU:71:LYS:HB2	1.67	0.60
25:BC:28:GLU:HB3	25:BC:29:PRO:HD3	1.84	0.60
23:DA:273(G):C:H2'	23:DA:274:G:H5''	1.84	0.60
12:CL:54:VAL:HG12	12:CL:55:ALA:H	1.65	0.60
23:BA:414:C:H2'	23:BA:415:A:C8	2.37	0.60
33:BK:53:LYS:HD2	33:BK:53:LYS:N	2.17	0.60
34:DL:38:GLN:HG3	34:DL:39:LYS:H	1.66	0.60
35:DM:76:LYS:N	35:DM:88:GLY:HA2	2.17	0.60
43:BU:8:LYS:NZ	43:BU:8:LYS:H	1.95	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
48:BZ:1:MET:SD	48:BZ:40:THR:HA	2.42	0.60
23:BA:1429:G:H2'	23:BA:1430:C:H6	1.66	0.60
46:BX:19:GLN:NE2	46:BX:41:ARG:HE	1.98	0.60
34:BL:71:VAL:HB	34:BL:72:PRO:HD3	1.84	0.60
23:BA:2776:A:H4'	23:BA:2777:G:H5''	1.83	0.60
18:AR:58:LEU:HB3	18:AR:62:GLU:HB2	1.83	0.60
23:DA:839:U:H2'	23:DA:840:C:C6	2.37	0.60
41:BS:40:ASN:O	41:BS:41:LYS:HG2	2.02	0.60
1:CA:299:G:H2'	1:CA:300:A:C8	2.36	0.60
12:CL:41:THR:OG1	12:CL:51:LEU:HB3	2.02	0.60
36:DN:11:ASN:OD1	36:DN:12:ARG:N	2.32	0.60
18:CR:58:LEU:HB3	18:CR:62:GLU:HB2	1.83	0.60
23:DA:153:C:OP1	46:DX:92:LYS:HE2	2.02	0.60
23:DA:2893:G:H5''	23:DA:2894:G:O4'	2.02	0.60
37:BO:24:LEU:HD13	37:BO:82:ILE:HG23	1.82	0.60
31:BI:4:LYS:HG2	31:BI:8:GLU:HG3	1.84	0.60
1:AA:1512:U:H2'	1:AA:1513:A:H8	1.65	0.60
23:DA:1603:A:C8	23:DA:1603:A:H5'	2.37	0.60
23:DA:1980:G:H3'	23:DA:1981:A:H5''	1.84	0.60
35:BM:30:GLY:HA2	35:BM:107:ALA:HB2	1.83	0.60
25:DC:28:GLU:HB3	25:DC:29:PRO:HD3	1.83	0.60
9:CI:97:LYS:HB3	9:CI:98:PRO:HD3	1.84	0.60
1:AA:1149:C:H2'	1:AA:1150:U:C6	2.37	0.60
23:BA:760:G:C2'	23:BA:761:A:H5'	2.31	0.59
1:CA:735:C:H2'	1:CA:736:C:H6	1.67	0.59
33:BK:3:GLN:CB	33:BK:4:PRO:HD2	2.29	0.59
35:BM:76:LYS:N	35:BM:88:GLY:HA2	2.17	0.59
23:BA:295:G:H4'	43:BU:2:ARG:NH1	2.16	0.59
23:BA:2038:G:H2'	23:BA:2039:C:C6	2.36	0.59
28:DF:83:ARG:HG3	28:DF:84:LYS:N	2.16	0.59
30:BH:113:ARG:HB2	30:BH:130:TYR:CE1	2.37	0.59
46:DX:19:GLN:NE2	46:DX:41:ARG:HE	2.00	0.59
35:BM:38:GLU:O	35:BM:127:ILE:HD13	2.02	0.59
1:AA:186(B):C:O2'	20:AT:89:ARG:HD2	2.02	0.59
23:BA:1803:A:H5''	23:BA:1804:C:OP2	2.01	0.59
41:BS:29:LEU:HD22	41:BS:69:LEU:HD11	1.83	0.59
23:DA:426:C:H2'	23:DA:427:U:H6	1.67	0.59
1:AA:841:U:O2'	1:AA:842:C:H5''	2.02	0.59
2:AB:95:GLN:HG3	2:AB:147:LYS:O	2.01	0.59
23:BA:2243:U:H2'	23:BA:2244:U:C6	2.37	0.59
23:DA:656:G:H2'	23:DA:657:U:O4'	2.02	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DA:1046:A:H1'	31:DI:4:LYS:HD3	1.83	0.59
12:AL:35:VAL:HG12	12:AL:36:CYS:O	2.01	0.59
2:CB:25:ASN:HD22	2:CB:25:ASN:N	1.99	0.59
23:DA:2038:G:H2'	23:DA:2039:C:C6	2.36	0.59
24:BB:70:C:H2'	24:BB:71:C:H6	1.68	0.59
23:DA:476:G:H4'	23:DA:502:A:N1	2.17	0.59
23:BA:2307:G:H2'	23:BA:2308:G:H5'	1.82	0.59
23:BA:2593:U:H2'	23:BA:2594:C:C6	2.37	0.59
27:BE:170:LEU:HD12	27:BE:171:PRO:HD2	1.84	0.59
1:AA:505:G:H2'	1:AA:506:G:H8	1.65	0.59
33:DK:53:LYS:N	33:DK:53:LYS:HD2	2.17	0.59
23:BA:1799:G:H8	25:BC:181:GLU:CD	2.06	0.59
44:BV:69:THR:HG22	44:BV:90:VAL:HG22	1.84	0.59
42:BT:47:PHE:HB3	42:BT:89:ILE:HD12	1.84	0.59
23:BA:826:U:O2	23:BA:832:G:C2	2.56	0.59
13:CM:75:ALA:O	13:CM:79:LYS:HG3	2.02	0.59
43:BU:76:CYS:CB	43:BU:77:PRO:HD2	2.32	0.59
11:AK:29:ILE:HG22	11:AK:44:SER:CB	2.31	0.59
23:BA:835:A:OP1	53:B5:52:LYS:HG2	2.02	0.59
1:AA:328:C:H4'	1:AA:329:A:C5'	2.30	0.59
26:DD:50:GLY:HA3	26:DD:75:VAL:HG11	1.83	0.59
24:BB:13:A:H5'	45:BW:74:ARG:NH2	2.16	0.59
26:BD:9:VAL:HG22	26:BD:25:VAL:HB	1.84	0.59
23:DA:2074:U:H2'	23:DA:2075:U:C6	2.38	0.59
23:DA:1540:G:H2'	23:DA:1541:U:O4'	2.02	0.59
1:AA:1187:G:H5'	9:AI:113:LYS:HE2	1.83	0.59
52:B4:21:ARG:HB3	52:B4:31:LEU:HD21	1.84	0.59
1:CA:1498:U:H1'	1:CA:1499:A:N7	2.17	0.59
7:CG:46:ALA:O	7:CG:50:ILE:HG12	2.02	0.59
39:BQ:104:GLN:HB3	40:BR:44:LYS:HZ1	1.67	0.59
23:DA:1264:G:OP1	50:D2:19:ARG:NH2	2.34	0.59
1:CA:1064:G:N2	1:CA:1190:G:H2'	2.18	0.59
1:CA:324:G:OP1	20:CT:70:SER:HB2	2.02	0.59
25:DC:108:PRO:HG3	25:DC:143:HIS:CE1	2.37	0.59
23:DA:942:G:H5'	34:DL:35:HIS:HB2	1.85	0.59
14:CN:32:SER:HB3	14:CN:41:ARG:HG2	1.83	0.59
47:DY:13:ALA:O	47:DY:17:SER:HA	2.02	0.59
35:BM:75:THR:HG21	35:BM:85:LYS:HZ2	1.66	0.59
23:DA:2712:U:H1'	23:DA:712(B):A:H8	1.66	0.59
36:BN:4:LEU:C	36:BN:6:SER:H	2.05	0.59
25:BC:267:SER:O	25:BC:270:ILE:HG13	2.02	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
12:CL:35:VAL:HG12	12:CL:36:CYS:O	2.01	0.59
3:CC:15:THR:HG21	3:CC:181:ASN:HA	1.84	0.59
23:BA:342:G:H2'	23:BA:343:C:H5''	1.84	0.59
3:AC:15:THR:HG21	3:AC:181:ASN:HA	1.83	0.59
17:AQ:64:PRO:HA	17:AQ:70:ARG:HG3	1.83	0.59
23:DA:185:U:H4'	23:DA:218:A:H4'	1.83	0.59
32:DJ:117:HIS:CE1	32:DJ:120:ARG:HE	2.19	0.59
1:CA:692:U:H5	11:CK:26:ASN:ND2	1.99	0.59
33:BK:104:ARG:NH1	33:BK:104:ARG:HB3	2.18	0.59
52:D4:21:ARG:HB3	52:D4:31:LEU:HD21	1.83	0.59
1:CA:805:C:H2'	1:CA:806:C:H6	1.67	0.59
1:AA:1104:G:H4'	2:AB:111:ARG:NH1	2.17	0.59
1:CA:304:U:H2'	1:CA:305:G:C8	2.37	0.59
1:CA:1111:A:H8	1:CA:1111:A:O5'	1.85	0.59
3:CC:112:SER:O	3:CC:116:VAL:HG23	2.02	0.59
40:DR:12:TYR:OH	40:DR:22:VAL:HG13	2.02	0.59
42:DT:12:VAL:HG12	42:DT:28:PHE:HA	1.84	0.59
32:DJ:39:ILE:HG22	32:DJ:40:ASP:O	2.03	0.59
1:AA:737:A:H2'	1:AA:738:C:C6	2.38	0.59
25:BC:44:ASN:HB3	25:BC:50:THR:HG21	1.84	0.59
35:DM:141:GLN:HG2	44:DV:72:ARG:HA	1.84	0.59
25:DC:35:LYS:HE3	25:DC:104:TYR:CD2	2.37	0.59
10:AJ:32:ALA:H	10:AJ:78:ASN:ND2	2.01	0.59
53:D5:57:ARG:HH11	53:D5:57:ARG:HB2	1.66	0.59
34:DL:57:THR:HG23	34:DL:59:LEU:HB2	1.85	0.59
36:DN:18:LEU:HD11	36:DN:22:ARG:CZ	2.32	0.59
19:AS:6:LYS:HD3	19:AS:7:LYS:HE2	1.83	0.59
26:BD:50:GLY:HA3	26:BD:75:VAL:HG11	1.84	0.59
46:BX:19:GLN:HE21	46:BX:41:ARG:HE	1.47	0.59
23:BA:661:C:O3'	34:BL:18:ARG:HG2	2.02	0.59
1:AA:1296:C:H5''	13:AM:44:ARG:HH22	1.68	0.59
23:BA:185:U:H2'	23:BA:186:G:H8	1.67	0.59
24:DB:87:G:H21	24:DB:89(B):A:H62	1.51	0.59
26:DD:173:VAL:HG12	26:DD:174:ASP:H	1.66	0.59
23:BA:839:U:H2'	23:BA:840:C:C6	2.37	0.59
1:CA:464:G:O6	1:CA:466:G:H5''	2.02	0.59
23:DA:2577:A:H5''	23:DA:2578:G:H5'	1.82	0.59
23:DA:1291:C:H2'	23:DA:1292:U:C6	2.37	0.59
2:AB:158:LEU:HD12	2:AB:158:LEU:H	1.68	0.59
13:CM:49:THR:HG22	13:CM:51:ALA:H	1.67	0.59
40:DR:81:TYR:C	40:DR:82:ARG:HG3	2.21	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:BL:91:PHE:CE2	34:BL:95:VAL:HG12	2.37	0.59
30:DH:92:VAL:HG21	30:DH:97:ILE:HD11	1.84	0.59
7:CG:102:ARG:HG2	7:CG:106:GLN:NE2	2.17	0.59
23:BA:330:A:H2	23:BA:1210:A:H2'	1.68	0.59
24:DB:13:A:H5'	45:DW:74:ARG:NH2	2.18	0.59
23:DA:779:U:OP1	25:DC:49:ILE:HG13	2.02	0.59
15:AO:5:LYS:HD3	15:AO:5:LYS:H	1.67	0.59
10:AJ:54:PHE:CD2	10:AJ:55:LYS:HG3	2.38	0.59
1:AA:1097:C:H2'	1:AA:1098:C:C6	2.37	0.59
23:BA:2476:A:H2'	23:BA:2476:A:N3	2.17	0.59
23:BA:442:G:H1'	27:BE:48:THR:HG21	1.85	0.59
23:BA:1264:G:OP1	50:B2:19:ARG:NH2	2.33	0.59
47:DY:38:GLN:HB3	47:DY:44:LEU:HB3	1.84	0.59
23:DA:1358:G:O2'	23:DA:1359:A:H5''	2.01	0.59
7:AG:26:PHE:O	7:AG:30:ILE:HG12	2.02	0.59
5:AE:47:LYS:HD3	5:AE:47:LYS:N	2.17	0.59
32:BJ:80:ALA:O	32:BJ:83:ILE:HG13	2.01	0.59
1:AA:1528:U:H6	1:AA:1528:U:H5''	1.67	0.59
32:DJ:119:GLU:O	32:DJ:123:GLU:HG3	2.02	0.59
40:BR:38:LEU:HD23	40:BR:39:LEU:N	2.17	0.59
23:BA:1794:U:H2'	23:BA:1795:C:C6	2.37	0.59
23:BA:2392:A:OP1	53:B5:32:LEU:HB3	2.02	0.59
23:BA:2496:C:OP1	35:BM:81:VAL:HG13	2.03	0.59
47:BY:13:ALA:O	47:BY:17:SER:HA	2.03	0.59
23:DA:295:G:H4'	43:DU:2:ARG:NH1	2.17	0.59
25:DC:30:GLU:CG	25:DC:63:ARG:HH21	2.16	0.59
12:CL:70:PRO:HD2	12:CL:101:ARG:HD3	1.82	0.59
27:DE:181:LEU:HD22	27:DE:186:ILE:HD11	1.85	0.59
1:AA:429:U:H1'	1:AA:430:A:H5''	1.85	0.59
39:DQ:50:ARG:NH2	40:DR:72:VAL:HG12	2.17	0.59
26:DD:36:ARG:HH12	26:DD:86:PRO:HD2	1.68	0.59
1:AA:33:A:H2'	1:AA:34:C:C6	2.38	0.59
23:DA:1799:G:H8	25:DC:181:GLU:CD	2.05	0.59
23:DA:1993:U:H4'	26:DD:128:SER:HB3	1.83	0.59
1:CA:190:G:H4'	1:CA:191(A):G:OP2	2.00	0.59
1:CA:828:A:H2'	1:CA:829:G:O4'	2.02	0.59
16:AP:13:HIS:C	16:AP:15:PRO:HD3	2.23	0.59
23:BA:779:U:H5'	25:BC:49:ILE:HD11	1.85	0.59
1:CA:841:U:O2'	1:CA:842:C:H5''	2.03	0.59
23:BA:83:G:H22	23:BA:102:G:H2'	1.65	0.59
11:CK:120:ARG:HH21	11:CK:126:ARG:HH21	1.49	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:946:A:H2'	1:CA:947:G:H8	1.66	0.59
40:BR:12:TYR:OH	40:BR:22:VAL:HG13	2.02	0.59
43:DU:13:VAL:HG11	43:DU:72:VAL:HB	1.83	0.59
28:DF:60:LEU:HD11	28:DF:92:VAL:CG1	2.31	0.59
52:B4:8:ASN:HD21	52:B4:11:LYS:H	1.51	0.59
23:DA:1311:G:C5'	23:DA:1311:G:H8	2.13	0.59
1:CA:552:U:O2	12:CL:30:PRO:HB3	2.03	0.59
39:BQ:50:ARG:NH2	40:BR:72:VAL:HG12	2.17	0.59
19:CS:63:THR:HG22	19:CS:66:MET:HG2	1.83	0.59
47:DY:38:GLN:O	47:DY:41:ILE:HG12	2.03	0.59
2:CB:95:GLN:HG3	2:CB:147:LYS:O	2.03	0.59
23:BA:814:C:C5	34:BL:27:HIS:NE2	2.71	0.59
11:AK:57:THR:HG22	11:AK:59:TYR:H	1.68	0.59
13:AM:49:THR:HG22	13:AM:51:ALA:H	1.66	0.59
27:BE:178:PRO:HB3	27:BE:201:VAL:HG11	1.83	0.59
1:CA:57:G:H2'	1:CA:58:C:C6	2.38	0.59
23:DA:1566:A:O2'	23:DA:1567:A:H5'	2.03	0.59
1:CA:114:U:H2'	1:CA:115:G:C8	2.37	0.59
23:BA:2346:A:H5''	23:BA:2383:G:H1'	1.84	0.59
23:BA:426:C:H2'	23:BA:427:U:H6	1.68	0.59
23:DA:2815:C:O2'	50:D2:43:HIS:HD2	1.86	0.59
40:BR:38:LEU:O	40:BR:39:LEU:HD13	2.03	0.59
42:BT:12:VAL:HG12	42:BT:28:PHE:HA	1.85	0.59
43:BU:13:VAL:HG11	43:BU:72:VAL:HB	1.84	0.59
49:B1:50:THR:HG22	49:B1:51:TYR:N	2.15	0.59
7:AG:102:ARG:HG2	7:AG:106:GLN:NE2	2.17	0.59
34:BL:57:THR:HG23	34:BL:59:LEU:CB	2.33	0.59
23:DA:674:G:C1'	27:DE:74:ARG:HD3	2.33	0.59
23:BA:1188:U:O2'	23:BA:1189:A:H5'	2.03	0.59
23:DA:2468:G:H22	23:DA:2481:G:H2'	1.67	0.59
30:BH:113:ARG:HB2	30:BH:130:TYR:CZ	2.38	0.59
13:CM:14:ARG:CZ	13:CM:42:ALA:HA	2.32	0.59
23:BA:2563:U:O2	23:BA:2565:A:H8	1.85	0.59
27:BE:192:LEU:HD21	27:BE:194:MET:HE3	1.84	0.59
39:BQ:36:ARG:HG2	39:BQ:40:PHE:CE1	2.38	0.59
23:DA:784:A:C5	25:DC:229:VAL:HG21	2.37	0.59
46:BX:58:ILE:HD11	46:BX:91:LYS:HG2	1.84	0.59
38:BP:132:LYS:O	38:BP:136:GLN:HG3	2.03	0.59
23:DA:826:U:O2	23:DA:832:G:C2	2.56	0.59
2:CB:158:LEU:HD12	2:CB:158:LEU:H	1.66	0.59
3:AC:91:LEU:HD13	3:AC:99:VAL:HB	1.84	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
27:DE:150:GLY:HA2	27:DE:172:TRP:CE3	2.37	0.59
3:CC:66:VAL:HB	3:CC:101:LEU:HD23	1.84	0.59
47:DY:2:LYS:HD2	47:DY:2:LYS:N	2.18	0.59
49:D1:50:THR:HG22	49:D1:51:TYR:N	2.15	0.59
1:AA:1391:U:H2'	1:AA:1392:G:H8	1.67	0.59
12:AL:35:VAL:HG22	12:AL:81:VAL:HG13	1.84	0.59
27:BE:80:ALA:O	27:BE:83:PHE:HB2	2.02	0.59
1:AA:464:G:O6	1:AA:466:G:H5'	2.03	0.59
3:CC:91:LEU:HD13	3:CC:99:VAL:HB	1.84	0.59
10:AJ:6:ILE:HG12	10:AJ:72:VAL:O	2.02	0.59
44:DV:14:LYS:HB2	44:DV:17:ALA:HB3	1.85	0.59
45:DW:21:LEU:H	45:DW:21:LEU:HD12	1.68	0.59
25:BC:95:LEU:O	25:BC:95:LEU:HD12	2.03	0.59
1:AA:1111:A:H8	1:AA:1111:A:O5'	1.86	0.59
44:BV:58:VAL:HA	44:BV:67:LEU:O	2.01	0.59
23:DA:2607:G:H2'	23:DA:2608:G:O4'	2.03	0.59
23:DA:760:G:C2'	23:DA:761:A:H5'	2.32	0.58
23:DA:835:A:OP1	53:D5:52:LYS:HG2	2.03	0.58
52:D4:34:ARG:HD2	52:D4:39:ARG:HG3	1.85	0.58
51:B3:11:LEU:HD13	51:B3:12:GLU:N	2.18	0.58
26:BD:33:VAL:HG12	26:BD:89:ASP:O	2.03	0.58
47:DY:24:LEU:HD22	47:DY:60:LEU:HD13	1.86	0.58
42:DT:47:PHE:HB3	42:DT:89:ILE:HD12	1.83	0.58
27:BE:150:GLY:HA2	27:BE:172:TRP:CE3	2.38	0.58
2:AB:118:LEU:HD13	2:AB:142:LEU:HA	1.84	0.58
38:DP:20:PRO:HD2	38:DP:86:ILE:HG23	1.84	0.58
38:DP:132:LYS:O	38:DP:136:GLN:HG3	2.02	0.58
37:BO:52:SER:HB2	37:BO:56:LEU:HB2	1.85	0.58
46:DX:37:ILE:CG2	46:DX:38:SER:N	2.66	0.58
40:BR:28:GLU:HB2	40:BR:31:ALA:HB2	1.85	0.58
33:DK:122:LEU:HD13	38:DP:72:VAL:HG11	1.83	0.58
7:AG:115:ARG:O	7:AG:118:VAL:HG22	2.02	0.58
4:AD:8:VAL:HB	4:AD:21:LEU:HD22	1.84	0.58
4:AD:108:LEU:HD21	4:AD:183:GLY:HA3	1.84	0.58
23:BA:1331:A:O2'	23:BA:1332:G:H8	1.86	0.58
25:DC:11:PRO:C	25:DC:13:ARG:H	2.07	0.58
13:CM:84:ILE:HD13	19:CS:66:MET:HE1	1.83	0.58
23:DA:414:C:H2'	23:DA:415:A:C8	2.38	0.58
44:BV:14:LYS:HB2	44:BV:17:ALA:HB3	1.85	0.58
1:AA:687:A:H4'	1:AA:688:G:O5'	2.04	0.58
4:CD:12:CYS:SG	4:CD:19:LEU:HB2	2.43	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
16:CP:13:HIS:C	16:CP:15:PRO:HD3	2.23	0.58
1:CA:1502:A:H5'	1:CA:1504:G:N7	2.18	0.58
40:DR:38:LEU:HD23	40:DR:39:LEU:N	2.18	0.58
23:BA:71:A:C2	42:BT:31:HIS:HE1	2.21	0.58
48:DZ:8:LEU:HB2	48:DZ:28:LEU:HD23	1.85	0.58
26:BD:132:HIS:CG	26:BD:135:HIS:NE2	2.72	0.58
23:BA:603:A:H61	23:BA:655:A:H1'	1.68	0.58
4:CD:108:LEU:HD21	4:CD:183:GLY:HA3	1.85	0.58
23:DA:1434:A:H2'	23:DA:1435:G:H8	1.68	0.58
23:DA:1478:G:O2'	23:DA:1558:A:H2	1.86	0.58
23:DA:2776:A:H4'	23:DA:2777:G:H5''	1.84	0.58
1:CA:892:A:H2'	1:CA:893:C:C6	2.38	0.58
1:AA:1064:G:N2	1:AA:1190:G:H2'	2.17	0.58
3:AC:89:GLU:O	3:AC:93:LYS:HB2	2.03	0.58
30:BH:53:ALA:O	30:BH:57:ARG:HB2	2.04	0.58
26:DD:1:MET:HB3	26:DD:83:ASP:O	2.02	0.58
2:AB:97:TRP:HZ2	2:AB:102:LEU:HD13	1.67	0.58
13:CM:27:LYS:HG3	13:CM:31:LYS:HE3	1.84	0.58
15:AO:56:LEU:HD21	23:BA:715:G:C2	2.38	0.58
1:AA:1498:U:H1'	1:AA:1499:A:N7	2.18	0.58
53:B5:37:SER:OG	53:B5:40:GLU:HG2	2.03	0.58
29:DG:13:LYS:HE2	29:DG:14:GLY:H	1.66	0.58
34:BL:32:THR:OG1	34:BL:36:LYS:HB3	2.04	0.58
35:BM:141:GLN:HG2	44:BV:72:ARG:HA	1.84	0.58
25:BC:271:ILE:O	25:BC:272:ALA:HB3	2.03	0.58
42:BT:63:LYS:HZ1	42:BT:72:LYS:HB3	1.67	0.58
23:DA:603:A:H61	23:DA:655:A:H1'	1.68	0.58
25:BC:30:GLU:CG	25:BC:63:ARG:HH21	2.16	0.58
4:CD:8:VAL:HB	4:CD:21:LEU:HD22	1.84	0.58
10:CJ:54:PHE:CD2	10:CJ:55:LYS:HG3	2.38	0.58
26:BD:36:ARG:HH12	26:BD:86:PRO:HD2	1.68	0.58
23:BA:7:G:H2'	23:BA:8:A:C8	2.37	0.58
15:AO:53:HIS:CE1	23:BA:715:G:O6	2.57	0.58
38:BP:20:PRO:HD2	38:BP:86:ILE:HG23	1.85	0.58
25:BC:70:TRP:CZ2	25:BC:150:LYS:HA	2.38	0.58
23:DA:2476:A:N3	23:DA:2476:A:H2'	2.17	0.58
6:AF:35:ALA:HA	6:AF:67:MET:HB3	1.85	0.58
23:BA:784:A:C5	25:BC:229:VAL:HG21	2.38	0.58
25:DC:8:PRO:HB3	25:DC:14:ARG:CB	2.33	0.58
23:BA:942:G:H5'	34:BL:35:HIS:HB2	1.86	0.58
23:DA:1161:C:O2'	40:DR:8:GLY:HA2	2.04	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
42:BT:26:TYR:O	42:BT:81:VAL:HG22	2.03	0.58
23:DA:886:C:C2'	23:DA:887:A:H4'	2.31	0.58
26:BD:92:THR:O	26:BD:95:ILE:HG13	2.03	0.58
36:DN:4:LEU:C	36:DN:6:SER:H	2.05	0.58
23:BA:1657:C:H2'	23:BA:1658:C:C6	2.38	0.58
1:CA:1327:C:H2'	1:CA:1328:C:C6	2.39	0.58
12:CL:35:VAL:HG22	12:CL:81:VAL:HG13	1.85	0.58
12:CL:82:VAL:HG21	12:CL:99:ILE:HD11	1.86	0.58
23:DA:528:A:C8	23:DA:528:A:C3'	2.87	0.58
30:DH:123:LEU:HD23	30:DH:124:GLY:N	2.18	0.58
27:BE:164:ARG:O	27:BE:168:ARG:HB2	2.03	0.58
23:BA:278:A:H4'	23:BA:279:C:OP1	2.04	0.58
23:BA:476:G:H4'	23:BA:502:A:N1	2.19	0.58
1:AA:304:U:H2'	1:AA:305:G:C8	2.39	0.58
23:BA:1966:A:H4'	23:BA:1967:C:OP1	2.02	0.58
32:BJ:119:GLU:O	32:BJ:123:GLU:HG3	2.04	0.58
23:DA:1952:A:C5	33:DK:22:ILE:HD12	2.39	0.58
41:DS:40:ASN:O	41:DS:41:LYS:HG2	2.04	0.58
2:CB:118:LEU:HD13	2:CB:142:LEU:HA	1.84	0.58
1:CA:1149:C:H2'	1:CA:1150:U:C6	2.37	0.58
20:AT:90:GLN:O	20:AT:93:GLU:HB3	2.03	0.58
39:DQ:95:LEU:O	39:DQ:98:LEU:HG	2.04	0.58
39:BQ:92:ARG:HG2	40:BR:11:GLN:NE2	2.19	0.58
32:BJ:39:ILE:HG22	32:BJ:40:ASP:O	2.03	0.58
23:DA:2392:A:OP1	53:D5:32:LEU:HB3	2.04	0.58
23:DA:71:A:C2	42:DT:31:HIS:HE1	2.22	0.58
26:DD:132:HIS:CG	26:DD:135:HIS:NE2	2.71	0.58
30:BH:114:LEU:HD21	30:BH:128:LEU:HD13	1.85	0.58
1:AA:707:C:H4'	11:AK:20:TYR:CD1	2.38	0.58
34:BL:18:ARG:HB3	34:BL:18:ARG:CZ	2.34	0.58
23:DA:558:G:OP1	32:DJ:134:PRO:HD2	2.03	0.58
25:DC:130:ALA:HB2	25:DC:192:THR:HB	1.86	0.58
23:DA:795:C:H2'	23:DA:796:C:C6	2.37	0.58
23:DA:7:G:H2'	23:DA:8:A:C8	2.37	0.58
23:DA:966:G:H2'	23:DA:967:C:H6	1.67	0.58
2:CB:97:TRP:HZ2	2:CB:102:LEU:HD13	1.69	0.58
47:BY:24:LEU:HD22	47:BY:60:LEU:HD13	1.85	0.58
30:DH:53:ALA:O	30:DH:57:ARG:HB2	2.03	0.58
9:AI:97:LYS:HB3	9:AI:98:PRO:HD3	1.84	0.58
1:AA:186(A):C:H5'	20:AT:78:ALA:HB1	1.85	0.58
23:BA:2393:A:H5''	34:BL:62:LEU:HB3	1.85	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
43:BU:45:VAL:HA	43:BU:62:GLU:HA	1.86	0.58
26:DD:92:THR:O	26:DD:95:ILE:HG13	2.03	0.58
3:CC:189:ALA:HB3	3:CC:196:LEU:HB3	1.86	0.58
24:DB:70:C:H2'	24:DB:71:C:H6	1.68	0.58
33:BK:47:ILE:HG13	33:BK:48:PRO:HD2	1.85	0.58
46:BX:27:GLU:HB3	46:BX:33:LYS:HG3	1.85	0.58
24:BB:11:C:H3'	24:BB:12:C:C6	2.39	0.58
10:AJ:6:ILE:HD11	10:AJ:72:VAL:HB	1.86	0.58
13:CM:27:LYS:HE2	13:CM:31:LYS:HE2	1.86	0.58
35:DM:66:ILE:HG22	35:DM:104:PHE:CD2	2.39	0.58
1:AA:1108:G:H5'	3:AC:176:HIS:CD2	2.38	0.58
23:BA:656:G:H2'	23:BA:657:U:O4'	2.02	0.58
23:DA:1378:A:O2'	23:DA:1379:A:H3'	2.03	0.58
23:BA:1603:A:H5'	23:BA:1603:A:C8	2.39	0.58
28:DF:36:LYS:HB3	28:DF:160:VAL:HB	1.84	0.58
34:DL:128:HIS:HA	34:DL:147:LEU:CB	2.21	0.58
7:CG:115:ARG:O	7:CG:118:VAL:HG22	2.03	0.58
23:BA:955:C:OP1	35:BM:85:LYS:HE2	2.04	0.58
47:BY:17:SER:HB3	47:BY:18:PRO:CD	2.30	0.58
23:DA:270(J):G:O2'	23:DA:270(K):G:H8	1.87	0.58
10:CJ:32:ALA:H	10:CJ:78:ASN:ND2	2.01	0.58
3:AC:47:LEU:HD21	3:AC:68:VAL:HG11	1.86	0.58
23:BA:1331:A:HO2'	23:BA:1332:G:H8	1.52	0.58
19:CS:5:LEU:HD12	19:CS:8:GLY:O	2.03	0.58
13:AM:84:ILE:HD13	19:AS:66:MET:HE1	1.86	0.58
10:CJ:16:LEU:HD12	10:CJ:70:ARG:HD2	1.86	0.58
28:DF:114:ILE:HG23	28:DF:115:ARG:HD2	1.84	0.58
1:AA:1327:C:H2'	1:AA:1328:C:C6	2.38	0.58
29:BG:13:LYS:HE2	29:BG:14:GLY:H	1.67	0.58
47:BY:38:GLN:O	47:BY:41:ILE:HG12	2.03	0.58
5:AE:137:GLU:O	5:AE:141:GLN:HG3	2.03	0.58
7:AG:115:ARG:O	7:AG:119:ARG:HG3	2.03	0.58
31:DI:4:LYS:HG2	31:DI:8:GLU:HG3	1.85	0.58
13:AM:75:ALA:O	13:AM:79:LYS:HG3	2.03	0.58
47:BY:2:LYS:HD2	47:BY:2:LYS:H	1.68	0.58
25:DC:31:LYS:O	25:DC:36:PRO:HD3	2.03	0.58
23:BA:1311:G:C8	23:BA:1311:G:C5'	2.86	0.58
28:BF:83:ARG:HG3	28:BF:84:LYS:N	2.17	0.58
23:DA:2039:C:H2'	23:DA:2040:C:C6	2.35	0.58
30:DH:113:ARG:HB2	30:DH:130:TYR:CE1	2.38	0.58
4:CD:173:TRP:CD1	4:CD:189:PRO:HG3	2.39	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:CC:35:GLU:O	3:CC:39:ILE:HG13	2.04	0.58
1:AA:1429:C:H2'	1:AA:1430:C:C6	2.39	0.58
29:DG:55:PRO:HG2	29:DG:61:HIS:CD2	2.38	0.58
23:BA:2294:C:H2'	23:BA:2295:C:C6	2.38	0.58
13:AM:27:LYS:HE2	13:AM:31:LYS:HE2	1.85	0.58
20:CT:90:GLN:O	20:CT:93:GLU:HB3	2.03	0.58
1:AA:828:A:H2'	1:AA:829:G:O4'	2.03	0.58
12:AL:37:THR:HG23	12:AL:38:VAL:H	1.69	0.58
37:DO:52:SER:HB2	37:DO:56:LEU:HB2	1.85	0.58
33:DK:86:ILE:H	33:DK:86:ILE:HD12	1.68	0.58
1:CA:250:A:H8	1:CA:250:A:O5'	1.87	0.58
23:BA:2393:A:C5'	34:BL:62:LEU:HD12	2.34	0.58
47:BY:2:LYS:HD2	47:BY:2:LYS:N	2.19	0.58
25:BC:35:LYS:HE3	25:BC:104:TYR:CD2	2.38	0.58
47:DY:16:LEU:HB2	47:DY:20:GLU:CG	2.34	0.58
27:DE:53:THR:HG22	27:DE:56:GLU:CD	2.23	0.58
30:BH:130:TYR:CD2	30:BH:132:PRO:HG3	2.39	0.58
44:DV:24:LEU:HD21	44:DV:86:VAL:CG2	2.33	0.58
20:CT:72:LEU:HD21	20:CT:76:ALA:C	2.24	0.58
10:AJ:13:HIS:HB3	10:AJ:68:HIS:NE2	2.19	0.58
19:AS:40:ILE:HD13	19:AS:62:ILE:HD11	1.86	0.58
53:D5:37:SER:OG	53:D5:40:GLU:HG2	2.04	0.58
1:AA:57:G:H2'	1:AA:58:C:C6	2.38	0.58
23:DA:814:C:C5	34:DL:27:HIS:NE2	2.72	0.58
1:CA:1320:C:H2'	1:CA:1321:C:O4'	2.04	0.58
8:AH:9:MET:HG3	8:AH:26:VAL:HG21	1.86	0.58
23:DA:114(B):A:H4'	32:DJ:48:ARG:HH22	1.69	0.57
36:DN:101:ALA:HB2	50:D2:44:THR:HG21	1.86	0.57
23:DA:245:G:O6	53:D5:8:LYS:HE3	2.04	0.57
23:DA:2439:A:C5'	23:DA:2439:A:C8	2.86	0.57
3:CC:22:TRP:HZ3	3:CC:24:ALA:HB2	1.68	0.57
23:BA:2468:G:H22	23:BA:2481:G:H2'	1.68	0.57
37:BO:26:LEU:O	37:BO:88:ASP:HB3	2.04	0.57
23:BA:1540:G:H2'	23:BA:1541:U:O4'	2.03	0.57
10:CJ:13:HIS:HB3	10:CJ:68:HIS:NE2	2.18	0.57
1:AA:1372:U:H2'	1:AA:1373:G:O4'	2.04	0.57
5:CE:48:ALA:HB2	5:CE:57:LYS:HD3	1.85	0.57
25:BC:11:PRO:C	25:BC:13:ARG:H	2.07	0.57
19:AS:63:THR:HG22	19:AS:66:MET:HG2	1.84	0.57
27:DE:80:ALA:O	27:DE:83:PHE:HB2	2.04	0.57
23:DA:83:G:N2	23:DA:102:G:H2'	2.18	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BA:83:G:N2	23:BA:102:G:H2'	2.19	0.57
1:CA:976:G:C8	1:CA:1358:U:H2'	2.39	0.57
40:DR:28:GLU:HB2	40:DR:31:ALA:HB2	1.85	0.57
23:DA:278:A:H4'	23:DA:279:C:OP1	2.04	0.57
1:AA:250:A:H8	1:AA:250:A:O5'	1.87	0.57
42:DT:57:LEU:N	42:DT:57:LEU:HD12	2.19	0.57
18:AR:54:ARG:HD2	18:AR:54:ARG:N	2.19	0.57
28:BF:47:LYS:HG3	28:BF:82:LEU:HD22	1.86	0.57
23:BA:1980:G:H3'	23:BA:1981:A:H5''	1.85	0.57
23:BA:832:G:OP1	34:BL:40:SER:HB3	2.03	0.57
23:BA:114(B):A:H4'	32:BJ:48:ARG:HH22	1.69	0.57
10:AJ:49:VAL:O	10:AJ:60:ARG:HB2	2.04	0.57
1:CA:979:C:H3'	1:CA:980:C:C5'	2.31	0.57
10:CJ:48:THR:HA	10:CJ:62:HIS:CB	2.29	0.57
2:CB:25:ASN:HB3	2:CB:27:LYS:HE2	1.86	0.57
23:BA:1568:G:P	25:BC:63:ARG:HH22	2.27	0.57
13:AM:44:ARG:HB2	13:AM:46:LYS:HG2	1.86	0.57
46:DX:83:GLU:HG2	46:DX:84:GLY:N	2.17	0.57
23:DA:1331:A:HO2'	23:DA:1332:G:H8	1.51	0.57
26:DD:84:PHE:CZ	26:DD:86:PRO:HG3	2.39	0.57
23:DA:1275:A:C8	36:DN:16:HIS:CD2	2.92	0.57
1:CA:1506:U:O2'	1:CA:1507:A:H5'	2.03	0.57
27:BE:195:ASP:OD2	27:BE:197:ASP:HB3	2.04	0.57
23:DA:1504:C:HO2'	23:DA:1505:C:P	2.26	0.57
23:DA:2633:G:O2'	26:DD:61:ARG:HD3	2.05	0.57
1:CA:262:A:H2'	1:CA:263:A:C8	2.39	0.57
5:CE:137:GLU:O	5:CE:141:GLN:HG3	2.03	0.57
23:BA:1566:A:O2'	23:BA:1567:A:H5'	2.04	0.57
23:DA:534:U:O2'	39:DQ:49:HIS:HD2	1.87	0.57
10:AJ:16:LEU:HD12	10:AJ:70:ARG:HD2	1.86	0.57
25:DC:246:PRO:HD2	25:DC:255:LYS:HD3	1.86	0.57
23:BA:2233:U:H2'	23:BA:2234:G:C8	2.40	0.57
28:DF:47:LYS:HG3	28:DF:82:LEU:HD22	1.85	0.57
11:CK:50:TYR:HB3	11:CK:54:ARG:HB2	1.86	0.57
1:CA:716:A:H1'	11:CK:118:GLY:HA2	1.85	0.57
8:CH:8:ASP:O	8:CH:12:ARG:HG2	2.05	0.57
23:BA:2647:U:H2'	23:BA:2648:C:C6	2.39	0.57
23:BA:1993:U:H4'	26:BD:128:SER:HB3	1.84	0.57
23:BA:2056:G:N2	23:BA:2057:A:H1'	2.20	0.57
2:AB:25:ASN:HB3	2:AB:27:LYS:HE2	1.85	0.57
1:AA:1320:C:H2'	1:AA:1321:C:O4'	2.04	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:DL:18:ARG:HB3	34:DL:18:ARG:CZ	2.33	0.57
23:BA:795:C:H2'	23:BA:796:C:C6	2.39	0.57
1:AA:1327:C:OP1	21:AU:20:LYS:HB3	2.04	0.57
23:BA:2104:G:H2'	23:BA:2105:C:C6	2.40	0.57
23:BA:1587:A:H2'	23:BA:1588:C:C6	2.39	0.57
30:BH:90:GLY:O	30:BH:91:SER:HB2	2.02	0.57
23:BA:1241:A:H5'	23:BA:1241:A:N3	2.19	0.57
23:DA:929:G:H8	23:DA:929:G:O5'	1.88	0.57
23:BA:221:A:H4'	23:BA:222:A:O5'	2.03	0.57
26:BD:1:MET:HB3	26:BD:83:ASP:O	2.03	0.57
23:DA:2243:U:H2'	23:DA:2244:U:C6	2.39	0.57
9:AI:3:GLN:HG2	9:AI:20:ARG:HG2	1.87	0.57
10:CJ:49:VAL:O	10:CJ:60:ARG:HB2	2.05	0.57
1:CA:1014:A:H4'	19:CS:14:HIS:CE1	2.40	0.57
26:BD:5:LEU:HB2	26:BD:51:PHE:CD2	2.39	0.57
23:BA:2376:A:N6	37:BO:89:ARG:HD2	2.20	0.57
51:D3:11:LEU:HD11	51:D3:51:GLU:HG3	1.87	0.57
1:CA:974:A:H8	1:CA:974:A:OP1	1.88	0.57
20:AT:72:LEU:HD21	20:AT:76:ALA:C	2.24	0.57
40:BR:81:TYR:C	40:BR:82:ARG:HG3	2.24	0.57
23:BA:46:C:OP2	23:BA:215:G:H2'	2.04	0.57
33:BK:86:ILE:H	33:BK:86:ILE:HD12	1.68	0.57
42:BT:57:LEU:HD12	42:BT:57:LEU:N	2.18	0.57
2:CB:162:ILE:HD12	2:CB:162:ILE:O	2.04	0.57
23:DA:572:A:C2	23:DA:2033:A:C2	2.93	0.57
34:BL:57:THR:HG23	34:BL:59:LEU:HB2	1.87	0.57
23:DA:1579:A:H5'	23:DA:1579:A:C8	2.35	0.57
26:DD:5:LEU:HB2	26:DD:51:PHE:CD2	2.39	0.57
1:AA:1294:G:H2'	1:AA:1295:G:C8	2.40	0.57
50:B2:33:CYS:SG	50:B2:40:LYS:HE3	2.45	0.57
51:D3:11:LEU:HD13	51:D3:12:GLU:N	2.19	0.57
1:AA:976:G:C8	1:AA:1358:U:H2'	2.39	0.57
35:BM:40:ALA:HB3	35:BM:127:ILE:HD11	1.87	0.57
27:DE:164:ARG:O	27:DE:168:ARG:HB2	2.05	0.57
36:BN:13:HIS:CE1	36:BN:16:HIS:HB2	2.39	0.57
23:BA:534:U:O2'	39:BQ:49:HIS:HD2	1.87	0.57
29:DG:20:ALA:HB1	29:DG:21:PRO:HD2	1.87	0.57
3:CC:89:GLU:O	3:CC:93:LYS:HB2	2.04	0.57
6:CF:35:ALA:HA	6:CF:67:MET:HB3	1.86	0.57
29:BG:121:ILE:HD11	29:BG:140:LYS:HD3	1.87	0.57
24:BB:30:C:OP2	37:BO:32:LEU:HD11	2.03	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DA:2647:U:H2'	23:DA:2648:C:C6	2.40	0.57
23:BA:2607:G:H2'	23:BA:2608:G:O4'	2.05	0.57
12:AL:116:ARG:HB3	12:AL:121:THR:O	2.04	0.57
8:AH:8:ASP:O	8:AH:12:ARG:HG2	2.04	0.57
24:DB:30:C:OP2	37:DO:32:LEU:HD11	2.04	0.57
46:DX:58:ILE:HD11	46:DX:91:LYS:HG2	1.85	0.57
32:BJ:148:GLY:HA3	32:BJ:149:PRO:O	2.05	0.57
9:CI:85:LEU:HD11	9:CI:96:LEU:HD22	1.86	0.57
23:DA:2790:A:H2'	23:DA:2791:C:H5''	1.87	0.57
51:D3:13:CYS:SG	51:D3:24:GLU:HG3	2.45	0.57
33:BK:122:LEU:HD13	38:BP:72:VAL:HG11	1.86	0.57
40:DR:64:HIS:CD2	40:DR:92:THR:HG22	2.39	0.57
43:DU:8:LYS:N	43:DU:8:LYS:NZ	2.51	0.57
2:AB:184:VAL:HG12	2:AB:197:VAL:HG13	1.87	0.57
53:D5:54:GLU:HA	53:D5:57:ARG:NH1	2.19	0.57
3:CC:22:TRP:CZ3	3:CC:24:ALA:HB2	2.40	0.57
30:BH:142:VAL:HG12	30:BH:143:SER:H	1.69	0.57
30:DH:113:ARG:HB2	30:DH:130:TYR:CZ	2.39	0.57
4:AD:134:ASP:O	4:AD:136:PRO:HD3	2.04	0.57
23:DA:1331:A:O2'	23:DA:1332:G:H8	1.88	0.57
27:BE:29:ASN:N	27:BE:112:MET:HE1	2.19	0.57
10:CJ:6:ILE:HD11	10:CJ:72:VAL:HB	1.85	0.57
27:DE:195:ASP:OD2	27:DE:197:ASP:HB3	2.05	0.57
38:BP:117:ASP:O	38:BP:121:ILE:HG13	2.05	0.57
23:DA:1154:G:H8	23:DA:1154:G:O5'	1.88	0.57
25:DC:231:HIS:CG	25:DC:232:PRO:HD2	2.39	0.57
23:DA:2104:G:H2'	23:DA:2105:C:C6	2.39	0.57
1:AA:222:U:H2'	1:AA:223:U:C6	2.39	0.57
24:BB:111:U:H2'	24:BB:112:G:H8	1.70	0.57
12:CL:46:LYS:HG2	12:CL:47:PRO:N	2.19	0.57
34:BL:135:LEU:O	34:BL:139:LYS:HB2	2.04	0.57
34:BL:114:ILE:HD12	34:BL:114:ILE:O	2.05	0.57
1:CA:1372:U:H2'	1:CA:1373:G:O4'	2.04	0.57
46:DX:27:GLU:CD	46:DX:33:LYS:HE3	2.24	0.57
1:CA:1070:U:H2'	1:CA:1071:C:H6	1.70	0.57
23:DA:832:G:OP1	34:DL:40:SER:HB3	2.05	0.57
1:CA:1097:C:H2'	1:CA:1098:C:C6	2.39	0.57
1:AA:1038:C:H2'	1:AA:1039:C:C6	2.39	0.57
28:DF:136:ARG:O	28:DF:154:GLY:HA2	2.05	0.57
14:AN:4:LYS:O	14:AN:7:ILE:HG13	2.05	0.57
1:CA:353:A:H8	1:CA:353:A:H5'	1.70	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
49:B1:48:ILE:H	49:B1:48:ILE:HD12	1.70	0.57
23:BA:966:G:H2'	23:BA:967:C:H6	1.69	0.57
1:CA:222:U:H2'	1:CA:223:U:C6	2.40	0.57
10:CJ:90:LEU:N	10:CJ:91:PRO:HD3	2.19	0.57
42:DT:26:TYR:O	42:DT:81:VAL:HG22	2.04	0.57
23:BA:1544:C:H3'	23:BA:1545:A:C5'	2.33	0.57
25:BC:217:ARG:NH1	25:BC:217:ARG:HG2	2.19	0.57
25:BC:130:ALA:HB2	25:BC:192:THR:HB	1.86	0.57
23:BA:1434:A:H2'	23:BA:1435:G:H8	1.67	0.57
1:CA:841:U:HO2'	1:CA:842:C:H6	1.51	0.57
47:BY:38:GLN:HB3	47:BY:44:LEU:HB3	1.84	0.57
1:CA:1320:C:H42	19:CS:36:ARG:HG3	1.69	0.57
23:BA:1819:A:H4'	23:BA:1820:U:H5''	1.86	0.57
29:BG:20:ALA:HB1	29:BG:21:PRO:HD2	1.87	0.57
23:DA:1360:A:H5'	23:DA:1361:G:OP2	2.05	0.57
1:CA:1294:G:H2'	1:CA:1295:G:C8	2.40	0.57
1:CA:1038:C:H2'	1:CA:1039:C:C6	2.40	0.57
47:BY:23:LYS:O	47:BY:27:GLU:HG3	2.04	0.57
40:BR:64:HIS:CD2	40:BR:92:THR:HG22	2.40	0.57
1:CA:684:A:H2'	1:CA:685:G:C8	2.39	0.57
25:BC:231:HIS:CG	25:BC:232:PRO:HD2	2.40	0.57
7:CG:115:ARG:O	7:CG:119:ARG:HG3	2.04	0.57
23:BA:1046:A:C3'	23:BA:1047:G:H5''	2.35	0.57
24:BB:43:C:H2'	24:BB:44:G:H5''	1.87	0.57
1:AA:1320:C:N4	19:AS:36:ARG:HG3	2.18	0.57
23:DA:1544:C:H3'	23:DA:1545:A:C5'	2.34	0.57
13:CM:44:ARG:HB2	13:CM:46:LYS:HG2	1.86	0.57
21:AU:9:ARG:O	21:AU:13:ILE:HD13	2.05	0.57
30:BH:6:LEU:HA	30:BH:15:VAL:HG13	1.87	0.57
29:BG:55:PRO:HG2	29:BG:61:HIS:CD2	2.40	0.57
25:DC:8:PRO:HB3	25:DC:14:ARG:HB3	1.87	0.57
23:BA:2746:U:H2'	23:BA:2747:G:H5'	1.87	0.57
8:CH:9:MET:HG3	8:CH:26:VAL:HG21	1.87	0.57
1:CA:237:C:H5''	17:CQ:25:ARG:CZ	2.35	0.57
25:BC:8:PRO:HB3	25:BC:14:ARG:CB	2.34	0.57
23:BA:529:A:H62	23:BA:2041:U:H3	1.53	0.57
4:AD:126:ILE:HG22	4:AD:127:THR:H	1.69	0.57
8:AH:102:ARG:N	8:AH:102:ARG:HE	2.03	0.57
1:AA:723:U:H5''	1:AA:724:G:OP2	2.05	0.57
42:DT:12:VAL:HG12	42:DT:27:THR:O	2.05	0.57
53:D5:53:PRO:HB2	53:D5:57:ARG:HH21	1.70	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
48:BZ:8:LEU:HB2	48:BZ:28:LEU:HD23	1.86	0.57
27:DE:52:LYS:HB3	27:DE:56:GLU:O	2.04	0.57
27:BE:52:LYS:HB3	27:BE:56:GLU:O	2.05	0.57
23:BA:528:A:C8	23:BA:528:A:C3'	2.88	0.57
26:BD:67:PHE:HE2	26:BD:75:VAL:HG22	1.70	0.57
26:BD:67:PHE:CE2	26:BD:75:VAL:HG22	2.40	0.57
12:CL:6:ILE:O	12:CL:10:VAL:HG23	2.05	0.57
26:DD:33:VAL:HG12	26:DD:89:ASP:O	2.04	0.57
23:DA:46:C:OP2	23:DA:215:G:H2'	2.05	0.57
1:AA:300:A:O5'	1:AA:300:A:H8	1.88	0.57
23:DA:319:C:H2'	23:DA:320:A:C8	2.40	0.57
44:BV:95:PRO:HB2	44:BV:127:LYS:HE3	1.87	0.57
1:CA:429:U:H1'	1:CA:430:A:H5''	1.85	0.57
1:CA:1075:C:H5''	2:CB:179:LYS:NZ	2.20	0.57
23:DA:581:C:OP1	39:DQ:31:SER:HB2	2.05	0.57
28:BF:136:ARG:O	28:BF:154:GLY:HA2	2.04	0.57
1:AA:663:A:H5''	18:AR:61:LYS:HE2	1.87	0.57
23:DA:2563:U:O2	23:DA:2565:A:H8	1.86	0.57
46:BX:37:ILE:CG2	46:BX:38:SER:N	2.68	0.57
23:BA:2633:G:O2'	26:BD:61:ARG:HD3	2.05	0.57
23:DA:301:G:C4	23:DA:302:C:C5	2.93	0.57
12:CL:116:ARG:HB3	12:CL:121:THR:O	2.05	0.57
2:AB:162:ILE:O	2:AB:162:ILE:HD12	2.05	0.56
23:BA:761:A:O5'	23:BA:761:A:C8	2.45	0.56
42:DT:31:HIS:ND1	42:DT:32:PRO:HD2	2.19	0.56
3:AC:22:TRP:HZ3	3:AC:24:ALA:HB2	1.70	0.56
23:DA:330:A:H2	23:DA:1210:A:H2'	1.70	0.56
23:BA:556:G:H2'	23:BA:557:U:H6	1.67	0.56
4:AD:26:CYS:HA	4:AD:31:CYS:HA	1.86	0.56
23:DA:2376:A:N6	37:DO:89:ARG:HD2	2.20	0.56
23:DA:2078:C:H2'	23:DA:2079:U:C6	2.38	0.56
23:BA:1275:A:C8	36:BN:16:HIS:CD2	2.93	0.56
27:BE:192:LEU:HD21	27:BE:194:MET:CE	2.35	0.56
23:BA:1360:A:H5'	23:BA:1361:G:OP2	2.05	0.56
4:CD:91:SER:HA	4:CD:94:LEU:HD12	1.87	0.56
34:DL:135:LEU:O	34:DL:139:LYS:HB2	2.04	0.56
1:AA:1016:A:H2'	1:AA:1017:G:O4'	2.05	0.56
23:BA:2790:A:H2'	23:BA:2791:C:H5''	1.87	0.56
1:CA:932:C:OP1	7:CG:4:ARG:HG2	2.04	0.56
8:AH:19:VAL:HG23	8:AH:21:LYS:HG2	1.87	0.56
1:AA:262:A:H2'	1:AA:263:A:C8	2.39	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
13:CM:113:PRO:O	13:CM:115:LYS:HD3	2.05	0.56
34:DL:111:ARG:HG3	34:DL:128:HIS:CB	2.36	0.56
28:BF:5:LEU:HD21	49:B1:50:THR:HG23	1.87	0.56
5:CE:33:VAL:HG11	5:CE:109:ILE:HD13	1.87	0.56
12:AL:6:ILE:O	12:AL:10:VAL:HG23	2.05	0.56
27:DE:203:GLN:HA	27:DE:206:ILE:O	2.05	0.56
23:BA:2892:A:H2'	23:BA:2893:G:H5'	1.87	0.56
38:BP:19:LEU:HD13	38:BP:78:LEU:HD22	1.87	0.56
28:BF:81:LYS:O	28:BF:82:LEU:HD23	2.04	0.56
1:CA:537:G:H5''	12:CL:112:ARG:NH2	2.20	0.56
47:DY:21:LEU:HD23	47:DY:22:GLU:N	2.19	0.56
39:DQ:36:ARG:HG2	39:DQ:40:PHE:CE1	2.40	0.56
27:DE:192:LEU:HD21	27:DE:194:MET:CE	2.36	0.56
23:BA:301:G:C4	23:BA:302:C:C5	2.93	0.56
24:BB:87:G:H21	24:BB:89(B):A:H62	1.52	0.56
24:DB:37:C:H2'	37:DO:95:HIS:HE1	1.70	0.56
39:DQ:68:ALA:O	39:DQ:71:GLN:HB3	2.04	0.56
25:BC:125:ILE:H	25:BC:125:ILE:HD12	1.70	0.56
1:CA:687:A:H4'	1:CA:688:G:O5'	2.05	0.56
27:BE:32:LEU:HD23	27:BE:32:LEU:C	2.26	0.56
3:CC:76:VAL:HG21	3:CC:103:VAL:HG11	1.88	0.56
23:DA:1495:A:H2'	23:DA:1496:A:N3	2.20	0.56
1:AA:237:C:H5''	17:AQ:25:ARG:CZ	2.34	0.56
23:BA:195:A:OP1	34:BL:46:LYS:HE2	2.06	0.56
23:DA:747:U:C4	50:D2:2:ALA:N	2.74	0.56
23:DA:806:C:O2'	23:DA:2445:G:H4'	2.05	0.56
25:DC:233:HIS:CE1	25:DC:247:ALA:H	2.23	0.56
23:DA:1046:A:C3'	23:DA:1047:G:H5''	2.35	0.56
28:DF:5:LEU:HD21	49:D1:50:THR:HG23	1.87	0.56
34:BL:115:LEU:HA	34:BL:134:ALA:CB	2.35	0.56
37:BO:49:VAL:HG13	37:BO:76:LYS:HD2	1.87	0.56
1:AA:979:C:H3'	1:AA:980:C:C5'	2.31	0.56
23:BA:2439:A:H8	23:BA:2439:A:H5'	1.68	0.56
3:CC:47:LEU:HD21	3:CC:68:VAL:HG11	1.86	0.56
23:BA:568:U:O4	40:BR:78:LYS:NZ	2.37	0.56
2:CB:88:ALA:HB2	2:CB:219:VAL:HG13	1.86	0.56
19:CS:29:ARG:HD3	19:CS:48:THR:HB	1.87	0.56
3:CC:7:PRO:HG3	3:CC:175:LEU:HD11	1.87	0.56
1:CA:1429:C:H2'	1:CA:1430:C:C6	2.40	0.56
14:CN:29:ARG:HG2	14:CN:31:ARG:O	2.05	0.56
4:AD:3:ARG:HD3	4:AD:5:ILE:CD1	2.36	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
44:DV:24:LEU:HD21	44:DV:86:VAL:HG21	1.87	0.56
33:DK:104:ARG:NH1	33:DK:104:ARG:HB3	2.20	0.56
34:DL:125:VAL:O	34:DL:145:PRO:HD2	2.05	0.56
32:BJ:62:ARG:HH21	32:BJ:64:ASP:HB2	1.69	0.56
42:DT:8:ILE:HD12	42:DT:8:ILE:N	2.20	0.56
10:AJ:6:ILE:HG22	10:AJ:98:ILE:HG23	1.88	0.56
23:BA:278:A:H2'	23:BA:279:C:O4'	2.06	0.56
23:DA:319:C:H2'	23:DA:320:A:H8	1.71	0.56
1:CA:1235:U:H5''	21:CU:3:LYS:HD2	1.87	0.56
23:BA:626:U:O2	34:BL:105:LEU:HB3	2.05	0.56
1:AA:353:A:H5'	1:AA:353:A:H8	1.69	0.56
4:AD:91:SER:HA	4:AD:94:LEU:HD12	1.87	0.56
13:AM:113:PRO:O	13:AM:115:LYS:HD3	2.06	0.56
1:AA:382:A:H2'	1:AA:383:A:C8	2.41	0.56
10:AJ:90:LEU:N	10:AJ:91:PRO:HD3	2.20	0.56
8:CH:102:ARG:HE	8:CH:102:ARG:N	2.04	0.56
47:DY:23:LYS:O	47:DY:27:GLU:HG3	2.05	0.56
43:BU:76:CYS:SG	43:BU:77:PRO:HD2	2.46	0.56
25:DC:44:ASN:HB3	25:DC:50:THR:CG2	2.36	0.56
7:AG:69:VAL:CA	7:AG:138:LYS:HD2	2.35	0.56
2:AB:88:ALA:HB2	2:AB:219:VAL:HG13	1.86	0.56
23:DA:1430:C:H2'	23:DA:1431:U:C6	2.40	0.56
19:AS:29:ARG:HD3	19:AS:48:THR:HB	1.88	0.56
51:B3:11:LEU:HD11	51:B3:51:GLU:HG3	1.86	0.56
23:BA:779:U:OP1	25:BC:49:ILE:HG13	2.04	0.56
23:DA:2305:A:H5''	28:DF:134:GLY:HA3	1.87	0.56
23:DA:1966:A:H4'	23:DA:1967:C:OP1	2.04	0.56
5:CE:127:ASN:O	5:CE:131:ILE:HG12	2.05	0.56
23:BA:1378:A:O2'	23:BA:1379:A:H3'	2.06	0.56
12:AL:54:VAL:HG12	12:AL:55:ALA:H	1.69	0.56
9:AI:85:LEU:HD11	9:AI:96:LEU:HD22	1.86	0.56
23:DA:1373:A:H2'	23:DA:1374:G:O4'	2.06	0.56
12:CL:37:THR:HG23	12:CL:38:VAL:H	1.70	0.56
23:BA:1952:A:C5	33:BK:22:ILE:HD12	2.41	0.56
1:AA:1502:A:H5'	1:AA:1504:G:N7	2.19	0.56
20:CT:53:LEU:O	20:CT:57:ARG:HD3	2.06	0.56
48:BZ:17:LYS:HD3	48:BZ:17:LYS:C	2.26	0.56
24:DB:8:U:H5''	37:DO:15:ARG:HH22	1.71	0.56
23:BA:828:U:H4'	23:BA:831:G:N1	2.20	0.56
34:BL:64:LYS:HD2	53:B5:25:MET:SD	2.46	0.56
47:DY:2:LYS:CD	47:DY:2:LYS:N	2.69	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
38:BP:28:VAL:HA	38:BP:89:VAL:HG12	1.86	0.56
1:AA:673:G:H5''	6:AF:87:ARG:HH12	1.71	0.56
6:CF:97:PHE:O	18:CR:31:LEU:HD23	2.05	0.56
3:AC:22:TRP:CZ3	3:AC:24:ALA:HB2	2.41	0.56
16:CP:20:VAL:HG21	16:CP:32:TYR:CG	2.40	0.56
26:DD:67:PHE:HE2	26:DD:75:VAL:HG22	1.70	0.56
4:CD:3:ARG:HD2	4:CD:3:ARG:N	2.21	0.56
2:AB:70:PHE:O	2:AB:71:VAL:HG13	2.05	0.56
26:BD:84:PHE:CZ	26:BD:86:PRO:HG3	2.41	0.56
34:DL:143:GLY:O	34:DL:145:PRO:HD3	2.05	0.56
23:DA:1187:G:H5''	40:DR:81:TYR:CE2	2.41	0.56
23:BA:1588:C:H2'	23:BA:1589:C:H6	1.70	0.56
1:CA:1117:G:H21	1:CA:1180:A:H1'	1.71	0.56
1:CA:555:C:H2'	1:CA:556:C:C6	2.41	0.56
1:CA:1191:A:H5''	3:CC:4:LYS:NZ	2.20	0.56
35:BM:66:ILE:HG22	35:BM:104:PHE:CD2	2.40	0.56
23:BA:706:A:H2'	23:BA:707:G:O4'	2.06	0.56
34:BL:128:HIS:HA	34:BL:147:LEU:CB	2.20	0.56
1:CA:520:A:OP2	12:CL:50:ALA:O	2.23	0.56
23:BA:886:C:C2'	23:BA:887:A:H4'	2.31	0.56
23:DA:2496:C:OP1	35:DM:81:VAL:HG13	2.05	0.56
23:DA:911:A:C6	35:DM:9:TYR:HE1	2.24	0.56
36:BN:51:LEU:HD22	36:BN:66:VAL:HG13	1.87	0.56
36:DN:51:LEU:HD22	36:DN:66:VAL:HG13	1.88	0.56
53:D5:14:VAL:HG13	53:D5:22:VAL:HG13	1.88	0.56
25:DC:182:LEU:O	25:DC:271:ILE:HD12	2.06	0.56
23:DA:1311:G:C5'	23:DA:1311:G:C8	2.87	0.56
12:AL:82:VAL:HG21	12:AL:99:ILE:HD11	1.86	0.56
24:DB:43:C:H2'	24:DB:44:G:H5''	1.86	0.56
36:BN:9:LYS:C	36:BN:10:LEU:HG	2.26	0.56
23:DA:1541:U:H3'	23:DA:1542:G:C3'	2.35	0.56
8:CH:51:VAL:HG21	8:CH:60:ARG:CG	2.35	0.56
23:DA:1434:A:H61	23:DA:1558:A:N6	2.03	0.56
6:AF:61:LEU:HB3	6:AF:63:TYR:HE2	1.71	0.56
10:CJ:6:ILE:HG22	10:CJ:98:ILE:HG23	1.87	0.56
15:AO:56:LEU:O	15:AO:60:VAL:HG23	2.06	0.56
4:CD:90:GLY:CA	4:CD:204:ILE:HD11	2.35	0.56
32:DJ:148:GLY:HA3	32:DJ:149:PRO:O	2.04	0.56
32:DJ:151:HIS:HD2	32:DJ:152:PRO:O	1.89	0.56
1:CA:1157:A:H61	1:CA:1178:G:H1'	1.71	0.56
44:DV:95:PRO:HB2	44:DV:127:LYS:HE3	1.88	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
38:BP:80:SER:HB3	38:BP:83:ILE:HG13	1.86	0.56
23:DA:2746:U:H2'	23:DA:2747:G:H5'	1.86	0.56
25:BC:126:GLN:O	25:BC:193:VAL:HG11	2.05	0.56
1:AA:116:A:H61	1:AA:313:A:H1'	1.70	0.56
3:AC:112:SER:O	3:AC:116:VAL:HG23	2.06	0.56
29:DG:121:ILE:HD11	29:DG:140:LYS:HD3	1.87	0.56
23:DA:1241:A:H5'	23:DA:1241:A:N3	2.21	0.56
17:AQ:14:LYS:HD2	17:AQ:14:LYS:H	1.71	0.56
13:CM:4:ILE:HA	13:CM:57:ARG:HG3	1.87	0.56
1:CA:109:A:C6	1:CA:326:G:C6	2.94	0.56
23:DA:2058:A:N6	23:DA:2059:A:N6	2.54	0.56
23:BA:1826:G:H4'	25:BC:242:ARG:NE	2.21	0.56
9:CI:3:GLN:HG2	9:CI:20:ARG:HG2	1.88	0.56
38:DP:28:VAL:HA	38:DP:89:VAL:HG12	1.88	0.56
23:BA:2744:G:N2	29:BG:143:GLN:HE22	2.01	0.56
3:AC:189:ALA:HB3	3:AC:196:LEU:HB3	1.86	0.56
27:BE:181:LEU:HD22	27:BE:186:ILE:HD11	1.88	0.56
32:BJ:90:LEU:O	32:BJ:111:GLU:HG3	2.06	0.56
39:BQ:58:ARG:O	39:BQ:62:ILE:HG12	2.05	0.56
52:D4:37:LYS:HD3	52:D4:39:ARG:HE	1.71	0.56
3:AC:77:ILE:C	3:AC:83:ARG:HB3	2.25	0.56
1:AA:946:A:H2'	1:AA:947:G:C8	2.40	0.56
35:DM:40:ALA:HB3	35:DM:127:ILE:HD11	1.87	0.56
36:DN:13:HIS:CE1	36:DN:16:HIS:HB2	2.39	0.56
23:BA:966:G:H2'	23:BA:967:C:C6	2.41	0.56
4:CD:126:ILE:HG22	4:CD:127:THR:H	1.70	0.56
1:CA:377:G:OP1	16:CP:3:LYS:HD2	2.06	0.56
1:AA:1070:U:H2'	1:AA:1071:C:H6	1.71	0.56
1:AA:160:A:H2'	1:AA:161:A:O4'	2.06	0.56
15:CO:28:GLN:O	15:CO:32:LEU:HG	2.06	0.56
1:AA:805:C:H2'	1:AA:806:C:H6	1.69	0.56
1:AA:59:A:H1'	1:AA:354:G:N2	2.20	0.56
27:DE:184:TYR:CE2	27:DE:188:ARG:HD2	2.40	0.56
23:DA:626:U:O2	34:DL:105:LEU:HB3	2.04	0.56
14:CN:4:LYS:O	14:CN:7:ILE:HG13	2.05	0.56
34:BL:33:ARG:O	34:BL:34:GLY:C	2.44	0.56
1:AA:942:G:N2	9:AI:124:GLN:HE22	2.02	0.56
23:DA:1005:C:O2'	32:DJ:51:THR:HG21	2.06	0.56
2:AB:27:LYS:CG	2:AB:194:PRO:HD2	2.36	0.56
52:B4:37:LYS:HD3	52:B4:39:ARG:HE	1.70	0.56
1:CA:1254:C:OP1	10:CJ:45:ARG:HD3	2.06	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
19:CS:40:ILE:HD13	19:CS:62:ILE:HD11	1.86	0.56
23:DA:828:U:H4'	23:DA:831:G:N1	2.21	0.56
2:AB:97:TRP:HH2	2:AB:176:GLU:CD	2.09	0.56
8:CH:19:VAL:HG23	8:CH:21:LYS:HG2	1.87	0.56
35:DM:58:PHE:O	35:DM:58:PHE:CD1	2.59	0.56
23:BA:1373:A:H2'	23:BA:1374:G:O4'	2.05	0.56
1:CA:819:A:H4'	1:CA:820:U:OP2	2.05	0.56
38:DP:117:ASP:O	38:DP:121:ILE:HG13	2.05	0.56
23:BA:581:C:H2'	23:BA:582:G:H8	1.69	0.56
18:CR:84:LYS:NZ	18:CR:84:LYS:HA	2.21	0.56
49:D1:48:ILE:H	49:D1:48:ILE:HD12	1.69	0.56
1:CA:99:C:C6	1:CA:99:C:H3'	2.41	0.56
7:AG:12:LEU:HD23	7:AG:12:LEU:H	1.71	0.56
23:DA:2233:U:H2'	23:DA:2234:G:C8	2.41	0.56
34:DL:33:ARG:O	34:DL:34:GLY:C	2.44	0.56
25:BC:233:HIS:CE1	25:BC:247:ALA:H	2.23	0.56
37:DO:35:ILE:CG1	37:DO:101:LEU:HD21	2.34	0.56
2:AB:80:ILE:HD11	2:AB:208:ILE:HG23	1.88	0.56
23:BA:572:A:C2	23:BA:2033:A:C2	2.93	0.56
23:BA:343:C:C5'	23:BA:343:C:H6	2.18	0.56
30:BH:12:LEU:H	30:BH:12:LEU:HD22	1.71	0.56
23:DA:2036:C:C6	23:DA:2036:C:H5'	2.38	0.56
37:DO:26:LEU:O	37:DO:88:ASP:HB3	2.06	0.56
40:DR:77:ALA:O	40:DR:79:VAL:N	2.39	0.56
8:AH:51:VAL:HG21	8:AH:60:ARG:CG	2.36	0.56
51:D3:11:LEU:HB2	51:D3:26:ASN:H	1.70	0.56
1:CA:255:G:H1'	17:CQ:16:GLN:NE2	2.21	0.56
21:CU:9:ARG:O	21:CU:13:ILE:HD13	2.05	0.56
23:DA:2294:C:H2'	23:DA:2295:C:C6	2.41	0.56
28:DF:110:ALA:O	28:DF:114:ILE:HG13	2.05	0.56
23:DA:999:U:H5''	23:DA:1154:G:O6	2.05	0.56
1:CA:1228:C:P	13:CM:108:ARG:HH22	2.27	0.56
1:AA:555:C:H2'	1:AA:556:C:C6	2.41	0.56
23:BA:276:A:H3'	23:BA:277:C:H5''	1.88	0.56
23:BA:1655:A:H1'	26:BD:113:PHE:CD2	2.41	0.56
10:AJ:38:ILE:HB	10:AJ:71:LEU:HB3	1.88	0.56
1:CA:382:A:H2'	1:CA:383:A:C8	2.40	0.56
1:CA:59:A:H1'	1:CA:354:G:N2	2.20	0.56
24:BB:8:U:H5''	37:BO:15:ARG:HH22	1.71	0.56
18:CR:54:ARG:N	18:CR:54:ARG:HD2	2.20	0.56
23:BA:1156:A:H4'	23:BA:1157:G:OP2	2.05	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1157:A:H61	1:AA:1178:G:H1'	1.71	0.56
43:DU:96:ILE:HG23	43:DU:101:LYS:O	2.06	0.56
50:D2:33:CYS:SG	50:D2:40:LYS:HE3	2.46	0.56
37:DO:49:VAL:HG13	37:DO:76:LYS:HD2	1.87	0.56
23:BA:1813:G:H1'	25:BC:50:THR:CG2	2.35	0.56
38:DP:48:ILE:H	38:DP:48:ILE:HD12	1.70	0.56
6:AF:97:PHE:O	18:AR:31:LEU:HD23	2.05	0.56
1:CA:977:A:C2'	1:CA:978:A:H5''	2.35	0.56
36:DN:9:LYS:C	36:DN:10:LEU:HG	2.26	0.56
19:CS:6:LYS:HD2	19:CS:6:LYS:H	1.71	0.56
3:AC:54:ARG:O	3:AC:69:HIS:HD2	1.89	0.56
4:AD:173:TRP:CD1	4:AD:189:PRO:HG3	2.41	0.56
3:CC:77:ILE:C	3:CC:83:ARG:HB3	2.26	0.56
28:BF:110:ALA:O	28:BF:114:ILE:HG13	2.06	0.56
25:DC:126:GLN:O	25:DC:193:VAL:HG11	2.05	0.56
23:DA:706:A:H2'	23:DA:707:G:O4'	2.07	0.56
41:DS:73:ALA:O	41:DS:106:ILE:HG12	2.05	0.56
51:B3:13:CYS:SG	51:B3:24:GLU:HG3	2.45	0.56
7:AG:38:LEU:O	7:AG:42:ILE:HG13	2.06	0.56
39:BQ:95:LEU:O	39:BQ:98:LEU:HG	2.05	0.55
42:BT:30:VAL:HG12	42:BT:31:HIS:N	2.21	0.55
25:BC:44:ASN:HB3	25:BC:50:THR:CG2	2.36	0.55
25:DC:271:ILE:O	25:DC:272:ALA:HB3	2.06	0.55
23:BA:2517:C:C6	23:BA:2542:A:C2	2.94	0.55
23:DA:2036:C:H6	23:DA:2036:C:C5'	2.18	0.55
35:BM:58:PHE:CD1	35:BM:58:PHE:O	2.59	0.55
23:DA:1996:C:OP1	33:DK:31:LYS:HE3	2.06	0.55
39:BQ:68:ALA:O	39:BQ:71:GLN:HB3	2.05	0.55
23:DA:442:G:H1'	27:DE:48:THR:HG21	1.88	0.55
16:AP:49:LEU:HD12	16:AP:50:LYS:H	1.71	0.55
24:DB:111:U:H2'	24:DB:112:G:H8	1.70	0.55
30:BH:92:VAL:HG21	30:BH:97:ILE:HD11	1.86	0.55
34:DL:114:ILE:HD12	34:DL:114:ILE:O	2.05	0.55
1:AA:1223:C:C5'	1:AA:1224:G:H5''	2.31	0.55
53:D5:52:LYS:N	53:D5:53:PRO:HD2	2.20	0.55
23:DA:2210:G:N2	23:DA:2211:G:H5'	2.21	0.55
23:BA:1173:G:H1'	23:BA:1177:A:H61	1.71	0.55
14:AN:29:ARG:HG2	14:AN:31:ARG:O	2.05	0.55
46:DX:27:GLU:CB	46:DX:33:LYS:HG3	2.37	0.55
4:CD:3:ARG:HD3	4:CD:5:ILE:CD1	2.37	0.55
32:DJ:62:ARG:HH21	32:DJ:64:ASP:HB2	1.70	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DA:1588:C:H2'	23:DA:1589:C:H6	1.71	0.55
1:AA:1530:G:H2'	1:AA:1531:A:C8	2.41	0.55
38:BP:59:THR:O	38:BP:78:LEU:HB2	2.06	0.55
25:BC:8:PRO:HB3	25:BC:14:ARG:HB3	1.88	0.55
6:CF:5:GLU:HB3	6:CF:62:TRP:HE1	1.71	0.55
1:AA:892:A:H2'	1:AA:893:C:C6	2.41	0.55
1:CA:918:A:H2'	1:CA:919:A:C8	2.41	0.55
1:CA:193:C:H2'	1:CA:194:C:C6	2.42	0.55
44:DV:8:TYR:HB2	44:DV:38:TYR:CZ	2.41	0.55
23:DA:221:A:H4'	23:DA:222:A:O5'	2.06	0.55
30:BH:92:VAL:HG23	30:BH:96:ASP:HB2	1.88	0.55
25:DC:242:ARG:CD	25:DC:242:ARG:H	2.08	0.55
33:BK:103:ALA:O	33:BK:106:LEU:HD13	2.06	0.55
23:BA:1430:C:H2'	23:BA:1431:U:C6	2.41	0.55
19:CS:6:LYS:HD3	19:CS:7:LYS:CE	2.36	0.55
34:DL:18:ARG:HB3	34:DL:18:ARG:NH1	2.22	0.55
28:DF:173:LEU:HD23	28:DF:176:LEU:HD12	1.88	0.55
36:BN:101:ALA:HB2	50:B2:44:THR:HG21	1.88	0.55
1:CA:405:U:H3'	1:CA:406:G:H5'	1.89	0.55
23:DA:64:A:O2'	42:DT:71:GLY:HA3	2.06	0.55
1:AA:7:G:H21	5:AE:121:LYS:CE	2.19	0.55
23:DA:278:A:H2'	23:DA:279:C:O4'	2.06	0.55
23:BA:1187:G:H5''	40:BR:81:TYR:CE2	2.41	0.55
1:CA:434:U:H2'	1:CA:435:C:C6	2.41	0.55
1:CA:723:U:H5''	1:CA:724:G:OP2	2.06	0.55
15:AO:24:SER:H	15:AO:27:VAL:HB	1.71	0.55
15:AO:28:GLN:O	15:AO:32:LEU:HG	2.07	0.55
39:DQ:104:GLN:HB3	40:DR:44:LYS:HZ1	1.71	0.55
38:DP:80:SER:HB3	38:DP:83:ILE:HG13	1.88	0.55
8:AH:50:ARG:H	8:AH:50:ARG:HD2	1.72	0.55
39:DQ:92:ARG:HG2	40:DR:11:GLN:NE2	2.21	0.55
39:BQ:92:ARG:HD2	39:BQ:95:LEU:CG	2.37	0.55
43:DU:45:VAL:HA	43:DU:62:GLU:HA	1.87	0.55
1:AA:691:G:O6	11:AK:52:GLY:HA2	2.06	0.55
12:AL:31:PHE:HB3	12:AL:83:LEU:HD22	1.89	0.55
23:DA:343:C:H6	23:DA:343:C:C5'	2.18	0.55
43:BU:71:LYS:HZ2	43:BU:71:LYS:HB2	1.70	0.55
40:BR:77:ALA:O	40:BR:79:VAL:N	2.39	0.55
23:BA:1541:U:H3'	23:BA:1542:G:C3'	2.37	0.55
51:B3:11:LEU:HB2	51:B3:26:ASN:H	1.70	0.55
4:AD:13:ARG:HB3	4:AD:38:TYR:O	2.07	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
27:BE:203:GLN:HA	27:BE:206:ILE:O	2.06	0.55
24:DB:11:C:H3'	24:DB:12:C:C6	2.41	0.55
38:DP:19:LEU:HD13	38:DP:78:LEU:HD22	1.89	0.55
1:AA:1506:U:O2'	1:AA:1507:A:H5'	2.06	0.55
48:DZ:11:SER:OG	48:DZ:13:ILE:HG12	2.06	0.55
1:AA:109:A:C6	1:AA:326:G:C6	2.95	0.55
6:AF:75:LEU:O	6:AF:79:LEU:HG	2.06	0.55
33:BK:73:ASP:OD1	33:BK:75:SER:HB3	2.07	0.55
12:AL:32:ARG:HE	12:AL:32:ARG:HA	1.72	0.55
23:BA:2058:A:N6	23:BA:2059:A:N6	2.54	0.55
23:BA:747:U:C4	50:B2:2:ALA:N	2.74	0.55
1:AA:1227:A:N3	1:AA:1227:A:H2'	2.22	0.55
34:DL:115:LEU:HA	34:DL:134:ALA:CB	2.36	0.55
47:DY:2:LYS:HA	47:DY:5:GLU:OE2	2.07	0.55
44:BV:76:LEU:N	44:BV:76:LEU:HD12	2.21	0.55
23:BA:1579:A:H5'	23:BA:1579:A:C8	2.33	0.55
1:AA:1320:C:N3	19:AS:72:GLY:HA3	2.21	0.55
30:BH:77:LEU:HD11	30:BH:101:LEU:HB2	1.88	0.55
46:BX:27:GLU:CD	46:BX:33:LYS:HE3	2.27	0.55
2:CB:70:PHE:O	2:CB:71:VAL:HG13	2.06	0.55
1:CA:376:G:OP2	16:CP:67:THR:HG21	2.07	0.55
23:BA:2749:A:H4'	29:BG:62:LYS:HB3	1.89	0.55
1:AA:1376:U:H2'	1:AA:1377:A:C8	2.42	0.55
44:BV:8:TYR:HB2	44:BV:38:TYR:CZ	2.41	0.55
20:AT:53:LEU:O	20:AT:57:ARG:HD3	2.07	0.55
12:AL:46:LYS:HG2	12:AL:47:PRO:N	2.21	0.55
47:BY:21:LEU:HD23	47:BY:22:GLU:N	2.22	0.55
1:AA:790:A:H5'	22:AV:6168:G:H4'	1.89	0.55
15:CO:24:SER:H	15:CO:27:VAL:HB	1.72	0.55
23:BA:195:A:H61	23:BA:198:C:H3'	1.72	0.55
23:DA:195:A:OP1	34:DL:46:LYS:HE2	2.07	0.55
43:DU:7:VAL:HB	43:DU:8:LYS:HZ2	1.71	0.55
23:BA:2393:A:H5'	34:BL:62:LEU:HD12	1.88	0.55
1:CA:1227:A:H2'	1:CA:1227:A:N3	2.21	0.55
23:DA:1046:A:H1'	31:DI:4:LYS:HD2	1.89	0.55
23:BA:2439:A:C5'	23:BA:2439:A:H8	2.20	0.55
23:BA:2210:G:N2	23:BA:2211:G:H5'	2.21	0.55
48:DZ:1:MET:HA	48:DZ:39:ASP:HB3	1.88	0.55
30:DH:12:LEU:H	30:DH:12:LEU:HD22	1.71	0.55
4:AD:3:ARG:HD2	4:AD:3:ARG:N	2.21	0.55
1:CA:406:G:H5''	4:CD:5:ILE:HD12	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
24:DB:104:A:O4'	44:DV:29:TYR:HE1	1.90	0.55
13:AM:106:ASN:O	13:AM:107:ALA:HB3	2.06	0.55
23:DA:2892:A:H2'	23:DA:2893:G:H5'	1.87	0.55
27:BE:192:LEU:HD23	27:BE:193:VAL:N	2.22	0.55
23:BA:2305:A:H5''	28:BF:134:GLY:HA3	1.88	0.55
34:BL:112:LEU:HD23	34:BL:113:LYS:N	2.22	0.55
45:DW:51:VAL:N	45:DW:62:LEU:HD12	2.22	0.55
23:DA:1639:U:H2'	23:DA:1640:C:H5''	1.89	0.55
11:AK:50:TYR:HB3	11:AK:54:ARG:HB2	1.87	0.55
1:AA:1004:A:H8	1:AA:1026:G:C8	2.25	0.55
36:DN:88:ARG:HG3	36:DN:89:ASP:OD1	2.07	0.55
23:DA:55:G:H2'	23:DA:56:A:H8	1.72	0.55
1:AA:255:G:H2'	1:AA:256:U:C6	2.42	0.55
23:BA:1996:C:OP1	33:BK:31:LYS:HE3	2.06	0.55
34:BL:111:ARG:HG3	34:BL:128:HIS:CB	2.36	0.55
32:DJ:42:GLU:HA	32:DJ:82:LYS:CB	2.32	0.55
3:CC:120:VAL:HG21	3:CC:137:ALA:HB2	1.89	0.55
52:D4:8:ASN:HD21	52:D4:11:LYS:H	1.52	0.55
23:DA:1173:G:H1'	23:DA:1177:A:H61	1.71	0.55
25:DC:71:ASP:HB3	25:DC:103:ARG:NH2	2.20	0.55
52:B4:34:ARG:HD2	52:B4:39:ARG:HG3	1.87	0.55
1:CA:1427:U:H2'	1:CA:1428:A:H8	1.72	0.55
3:AC:35:GLU:O	3:AC:39:ILE:HG13	2.05	0.55
1:CA:255:G:H2'	1:CA:256:U:C6	2.42	0.55
35:DM:112:GLU:CD	35:DM:112:GLU:H	2.10	0.55
23:BA:1292:U:H2'	23:BA:1293:C:C6	2.42	0.55
1:CA:1530:G:H2'	1:CA:1531:A:C8	2.41	0.55
38:DP:57:PHE:O	38:DP:59:THR:N	2.39	0.55
46:DX:23:LYS:HB3	46:DX:37:ILE:HG12	1.87	0.55
1:AA:601:C:H2'	1:AA:602:A:H8	1.72	0.55
1:AA:819:A:H4'	1:AA:820:U:OP2	2.05	0.55
1:AA:99:C:C6	1:AA:99:C:H3'	2.42	0.55
23:BA:2815:C:O2'	50:B2:43:HIS:HD2	1.88	0.55
13:AM:82:MET:HG3	23:BA:888:C:H5'	1.89	0.55
29:BG:44:VAL:O	29:BG:50:VAL:HG13	2.07	0.55
1:AA:1279:A:H62	3:AC:26:LYS:HE2	1.71	0.55
23:BA:1971:A:C4	25:BC:241:PRO:HG3	2.41	0.55
42:BT:12:VAL:HG12	42:BT:27:THR:O	2.07	0.55
1:AA:1224:G:H4'	13:AM:102:ARG:NH2	2.21	0.55
17:AQ:82:MET:O	17:AQ:86:GLU:HG2	2.06	0.55
2:CB:27:LYS:CG	2:CB:194:PRO:HD2	2.36	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:DD:67:PHE:CE2	26:DD:75:VAL:HG22	2.41	0.55
19:AS:6:LYS:HD3	19:AS:7:LYS:CE	2.36	0.55
23:BA:2036:C:H5'	23:BA:2036:C:C6	2.39	0.55
14:CN:24:CYS:HB3	14:CN:29:ARG:N	2.20	0.55
11:AK:120:ARG:HH21	11:AK:126:ARG:NH2	2.05	0.55
23:DA:1292:U:H2'	23:DA:1293:C:C6	2.42	0.55
23:BA:775:G:C4	23:BA:794:G:C8	2.94	0.55
2:CB:97:TRP:HH2	2:CB:176:GLU:CD	2.09	0.55
23:BA:2648:C:H2'	23:BA:2649:U:C6	2.42	0.55
46:BX:23:LYS:HB3	46:BX:37:ILE:HG12	1.89	0.55
1:AA:1501:C:C4	1:AA:1504:G:C6	2.95	0.55
29:DG:44:VAL:O	29:DG:50:VAL:HG13	2.07	0.55
12:AL:41:THR:OG1	12:AL:51:LEU:HB3	2.06	0.55
23:BA:483:A:H4'	43:BU:49:VAL:HG23	1.89	0.55
1:CA:1016:A:H2'	1:CA:1017:G:O4'	2.06	0.55
3:AC:76:VAL:HG21	3:AC:103:VAL:HG11	1.88	0.55
43:BU:59:GLY:C	43:BU:61:ILE:H	2.10	0.55
1:AA:1117:G:H21	1:AA:1180:A:H1'	1.70	0.55
34:DL:33:ARG:HG3	34:DL:36:LYS:CD	2.22	0.55
23:BA:2014:A:H2'	23:BA:2015:A:C8	2.42	0.55
23:DA:1826:G:H4'	25:DC:242:ARG:NE	2.22	0.55
2:CB:184:VAL:HG12	2:CB:197:VAL:HG13	1.89	0.55
23:DA:676:A:H2	23:DA:802:A:H61	1.54	0.55
1:AA:692:U:H5	11:AK:26:ASN:ND2	2.03	0.55
53:B5:14:VAL:HG13	53:B5:22:VAL:HG13	1.89	0.55
16:AP:20:VAL:HG21	16:AP:32:TYR:CG	2.42	0.55
25:BC:182:LEU:O	25:BC:271:ILE:HD12	2.06	0.55
3:CC:11:ARG:HB3	3:CC:15:THR:HB	1.89	0.55
10:CJ:8:LEU:HG	10:CJ:96:ILE:HG22	1.89	0.55
19:AS:6:LYS:HD2	19:AS:6:LYS:H	1.71	0.55
48:BZ:1:MET:HA	48:BZ:39:ASP:HB3	1.88	0.55
23:DA:1188:U:O2'	23:DA:1189:A:H5'	2.06	0.55
4:CD:134:ASP:O	4:CD:136:PRO:HD3	2.06	0.55
23:BA:323:G:H5'	27:BE:169:ASN:HD21	1.71	0.55
23:DA:1276:A:O2'	36:DN:16:HIS:HE1	1.90	0.55
15:AO:53:HIS:HE1	23:BA:715:G:O6	1.90	0.55
1:CA:976:G:H8	1:CA:1358:U:H2'	1.72	0.55
28:DF:81:LYS:O	28:DF:82:LEU:HD23	2.06	0.55
1:AA:1504:G:OP1	1:AA:1507:A:H4'	2.07	0.55
7:CG:38:LEU:O	7:CG:42:ILE:HG13	2.07	0.55
4:AD:155:LEU:O	4:AD:159:ARG:HG2	2.06	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
33:BK:24:VAL:HG23	33:BK:33:ALA:HB2	1.89	0.55
1:CA:1376:U:H2'	1:CA:1377:A:C8	2.41	0.55
1:AA:149:A:H2'	1:AA:150:C:C6	2.42	0.55
8:CH:50:ARG:HD2	8:CH:50:ARG:H	1.72	0.55
26:BD:111:ARG:HD2	26:BD:160:TYR:CE1	2.42	0.55
1:CA:1137:C:H4'	1:CA:1138:G:C2	2.42	0.55
35:DM:75:THR:HA	35:DM:88:GLY:HA3	1.88	0.55
17:AQ:86:GLU:O	17:AQ:90:ILE:HG12	2.07	0.55
53:B5:52:LYS:N	53:B5:53:PRO:HD2	2.22	0.55
48:DZ:26:LEU:HB2	48:DZ:28:LEU:HD13	1.89	0.55
23:BA:2543:G:C8	23:BA:2543:G:H5'	2.40	0.55
10:AJ:8:LEU:HG	10:AJ:96:ILE:HG22	1.88	0.55
30:BH:116:LEU:HD22	30:BH:128:LEU:HD21	1.89	0.55
35:DM:45:GLN:CD	35:DM:45:GLN:H	2.10	0.55
23:DA:1448:G:H2'	23:DA:149(B):A:C8	2.41	0.55
1:CA:946:A:H2'	1:CA:947:G:C8	2.41	0.55
1:CA:18:C:H5''	5:CE:127:ASN:ND2	2.22	0.55
2:CB:235:SER:O	2:CB:239:VAL:HG23	2.06	0.55
2:AB:235:SER:O	2:AB:239:VAL:HG23	2.07	0.55
4:CD:13:ARG:HB3	4:CD:38:TYR:O	2.07	0.55
41:BS:4:LYS:HG2	41:BS:106:ILE:HG22	1.88	0.55
23:DA:1833:U:H2'	23:DA:1834:U:H6	1.70	0.55
4:AD:90:GLY:CA	4:AD:204:ILE:HD11	2.36	0.55
12:AL:53:LYS:HD2	12:AL:53:LYS:N	2.22	0.55
1:CA:971:G:H1'	1:CA:1365:G:O2'	2.07	0.55
25:DC:217:ARG:HG2	25:DC:217:ARG:NH1	2.19	0.54
1:CA:300:A:H8	1:CA:300:A:O5'	1.90	0.54
36:DN:12:ARG:HD3	36:DN:16:HIS:ND1	2.23	0.54
1:CA:1065:U:H4'	1:CA:1066:C:O5'	2.07	0.54
1:CA:624:C:O3'	16:CP:10:GLY:HA2	2.07	0.54
6:CF:75:LEU:O	6:CF:79:LEU:HG	2.07	0.54
23:DA:2102:U:H2'	23:DA:2103:C:C6	2.43	0.54
1:AA:1137:C:H4'	1:AA:1138:G:C2	2.42	0.54
23:BA:380:U:O2'	46:BX:20:ARG:HB3	2.08	0.54
12:CL:53:LYS:N	12:CL:53:LYS:HD2	2.22	0.54
8:AH:58:TYR:O	8:AH:59:LEU:HD23	2.07	0.54
41:BS:103:ILE:HD12	41:BS:103:ILE:H	1.73	0.54
42:DT:28:PHE:HE2	42:DT:92:LEU:HD11	1.71	0.54
23:DA:2393:A:H5'	34:DL:62:LEU:HD12	1.89	0.54
34:BL:84:ASN:HB3	34:BL:86:LYS:HG2	1.90	0.54
49:B1:46:ASN:HB2	49:B1:64:LYS:HB2	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DM:140:ALA:HB3	44:DV:53:ILE:HD13	1.89	0.54
48:BZ:26:LEU:HB2	48:BZ:28:LEU:HD13	1.89	0.54
23:BA:1005:C:O2'	32:BJ:51:THR:HG21	2.06	0.54
23:BA:1510:A:H2'	23:BA:1511:A:H8	1.72	0.54
33:DK:68:GLU:HB3	33:DK:78:ARG:HB2	1.90	0.54
37:BO:34:HIS:ND1	37:BO:54:LEU:HB2	2.22	0.54
32:DJ:90:LEU:O	32:DJ:111:GLU:HG3	2.07	0.54
30:DH:114:LEU:HD21	30:DH:128:LEU:HD13	1.89	0.54
23:DA:1614:A:H62	41:DS:93:ALA:HB2	1.72	0.54
41:DS:4:LYS:HG2	41:DS:106:ILE:HG22	1.89	0.54
23:BA:270(Q):C:HO2'	23:BA:270(R):C:H6	1.50	0.54
35:DM:26:TYR:HA	44:DV:81:ARG:HH21	1.72	0.54
23:BA:1932:A:H3'	23:BA:1933:G:H8	1.72	0.54
23:BA:1266:G:H5''	50:B2:23:HIS:NE2	2.21	0.54
23:BA:806:C:O2'	23:BA:2445:G:H4'	2.07	0.54
35:BM:81:VAL:CG1	35:BM:82:ARG:HG2	2.37	0.54
25:DC:25:THR:HG21	25:DC:81:ALA:HB1	1.89	0.54
1:CA:1347:G:H8	9:CI:107:ARG:HB3	1.70	0.54
28:BF:173:LEU:HD23	28:BF:176:LEU:HD12	1.89	0.54
23:BA:2036:C:C5'	23:BA:2036:C:H6	2.20	0.54
23:BA:1448:G:H2'	23:BA:149(B):A:C8	2.42	0.54
1:CA:1152:A:H2'	1:CA:1153:C:H6	1.73	0.54
13:AM:34:LEU:HD22	13:AM:39:ILE:HB	1.89	0.54
6:CF:12:PRO:HG2	6:CF:55:ASP:OD2	2.07	0.54
23:DA:1257:C:H4'	27:DE:83:PHE:CE2	2.42	0.54
1:CA:1501:C:C4	1:CA:1504:G:C6	2.94	0.54
27:DE:192:LEU:HD23	27:DE:193:VAL:N	2.22	0.54
1:AA:1157:A:H4'	1:AA:1158:C:O5'	2.08	0.54
1:AA:1065:U:H4'	1:AA:1066:C:O5'	2.07	0.54
1:CA:363:A:C8	12:CL:32:ARG:NH2	2.76	0.54
1:CA:269:C:H2'	1:CA:270:A:C8	2.42	0.54
35:BM:26:TYR:HA	44:BV:81:ARG:HH21	1.72	0.54
1:AA:974:A:OP1	1:AA:974:A:H8	1.89	0.54
7:CG:12:LEU:H	7:CG:12:LEU:HD23	1.71	0.54
27:DE:32:LEU:C	27:DE:32:LEU:HD23	2.28	0.54
23:BA:1647:G:H3'	23:BA:1647:G:OP2	2.08	0.54
23:BA:999:U:H5''	23:BA:1154:G:O6	2.06	0.54
33:DK:73:ASP:OD1	33:DK:75:SER:HB3	2.07	0.54
23:DA:2014:A:H2'	23:DA:2015:A:C8	2.43	0.54
23:BA:1046:A:H1'	31:BI:4:LYS:HD2	1.89	0.54
23:DA:910:A:C4	35:DM:13:GLN:NE2	2.74	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
21:AU:18:TYR:O	21:AU:22:ARG:HB3	2.07	0.54
35:BM:8:LYS:O	35:BM:9:TYR:HB3	2.08	0.54
26:BD:101:ARG:HD3	26:BD:169:ASN:ND2	2.22	0.54
3:CC:35:GLU:HA	3:CC:38:ARG:HG2	1.90	0.54
4:AD:108:LEU:HD23	4:AD:110:PHE:CE2	2.42	0.54
3:CC:54:ARG:O	3:CC:69:HIS:HD2	1.90	0.54
9:CI:17:VAL:HG13	9:CI:63:ILE:HD11	1.90	0.54
23:BA:1434:A:H61	23:BA:1558:A:N6	2.05	0.54
34:BL:143:GLY:O	34:BL:145:PRO:HD3	2.08	0.54
38:DP:59:THR:O	38:DP:78:LEU:HB2	2.07	0.54
8:AH:97:VAL:HG13	8:AH:98:LYS:N	2.23	0.54
1:CA:1100:C:OP2	2:CB:96:ARG:HG2	2.08	0.54
34:DL:112:LEU:HD23	34:DL:113:LYS:N	2.22	0.54
13:AM:15:VAL:O	13:AM:19:LEU:HD23	2.08	0.54
13:CM:10:PRO:HB2	13:CM:18:ALA:HB1	1.90	0.54
23:BA:1639:U:H2'	23:BA:1640:C:H5''	1.88	0.54
17:CQ:14:LYS:HD2	17:CQ:14:LYS:H	1.72	0.54
18:AR:84:LYS:HA	18:AR:84:LYS:NZ	2.22	0.54
23:BA:137(B):G:H2'	23:BA:139:G:N7	2.23	0.54
27:DE:183:VAL:O	27:DE:187:VAL:HG23	2.08	0.54
40:DR:6:LYS:O	40:DR:37:VAL:HG21	2.08	0.54
30:DH:92:VAL:HG23	30:DH:96:ASP:HB2	1.90	0.54
34:DL:64:LYS:HD2	53:D5:25:MET:SD	2.48	0.54
21:CU:18:TYR:O	21:CU:22:ARG:HB3	2.08	0.54
23:BA:270(J):G:O2'	23:BA:270(K):G:H8	1.87	0.54
23:BA:910:A:C4	35:BM:13:GLN:NE2	2.76	0.54
45:BW:35:ASN:N	45:BW:35:ASN:ND2	2.55	0.54
1:AA:977:A:C2'	1:AA:978:A:H5''	2.35	0.54
52:B4:34:ARG:HB3	52:B4:42:LEU:HD22	1.88	0.54
3:AC:7:PRO:HG3	3:AC:175:LEU:HD11	1.89	0.54
23:DA:819:A:C4	23:DA:1189:A:C2	2.95	0.54
23:BA:1614:A:H61	41:BS:88:ARG:H	1.56	0.54
1:AA:1152:A:H2'	1:AA:1153:C:H6	1.73	0.54
34:BL:125:VAL:O	34:BL:145:PRO:HD2	2.07	0.54
13:CM:106:ASN:O	13:CM:107:ALA:HB3	2.06	0.54
1:AA:793:U:H3'	1:AA:794:A:C5'	2.36	0.54
23:DA:966:G:H2'	23:DA:967:C:C6	2.41	0.54
1:AA:1017:G:H2'	1:AA:1018:C:C6	2.43	0.54
1:CA:1017:G:H2'	1:CA:1018:C:C6	2.42	0.54
33:DK:96:THR:O	33:DK:97:ARG:C	2.46	0.54
23:DA:2749:A:H4'	29:DG:62:LYS:HB3	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1004:A:H8	1:CA:1026:G:C8	2.25	0.54
1:CA:559:A:H4'	1:CA:560:U:H5''	1.90	0.54
23:BA:1495:A:H2'	23:BA:1496:A:N3	2.22	0.54
23:DA:919:G:H5'	24:DB:81:G:H1'	1.89	0.54
3:AC:120:VAL:HG21	3:AC:137:ALA:HB2	1.88	0.54
47:DY:11:GLU:OE1	47:DY:11:GLU:N	2.40	0.54
29:DG:27:LYS:HG2	29:DG:32:GLU:HB2	1.90	0.54
1:CA:1223:C:P	19:CS:78:ARG:HH21	2.29	0.54
35:DM:141:GLN:HA	44:DV:71:VAL:O	2.08	0.54
44:DV:74:VAL:O	44:DV:76:LEU:HD12	2.07	0.54
23:BA:948:G:H8	23:BA:948:G:C5'	2.20	0.54
27:DE:63:LYS:HE3	27:DE:75:HIS:O	2.08	0.54
33:DK:103:ALA:O	33:DK:106:LEU:HD13	2.07	0.54
4:CD:26:CYS:HA	4:CD:31:CYS:HA	1.90	0.54
1:CA:692:U:H5	11:CK:26:ASN:HD22	1.55	0.54
23:BA:64:A:O2'	42:BT:71:GLY:HA3	2.07	0.54
1:CA:793:U:H3'	1:CA:794:A:C5'	2.37	0.54
1:AA:176:C:H5''	20:AT:29:LYS:NZ	2.23	0.54
20:AT:97:ALA:O	20:AT:99:LEU:N	2.41	0.54
36:BN:88:ARG:HG3	36:BN:89:ASP:OD1	2.06	0.54
23:BA:1833:U:H2'	23:BA:1834:U:H6	1.72	0.54
23:BA:576:U:H2'	23:BA:577:G:C8	2.43	0.54
23:BA:319:C:H2'	23:BA:320:A:C8	2.42	0.54
23:DA:2774:C:H2'	23:DA:2775:A:O4'	2.06	0.54
23:BA:1161:C:O2'	40:BR:23:GLU:HG2	2.08	0.54
40:BR:38:LEU:HD22	40:BR:52:VAL:HG11	1.90	0.54
42:BT:23:GLU:HG3	42:BT:24:GLY:H	1.72	0.54
5:CE:76:ILE:HD11	5:CE:142:LEU:HD11	1.90	0.54
23:DA:2744:G:H21	29:DG:143:GLN:NE2	2.00	0.54
17:AQ:56:VAL:HG23	17:AQ:81:ARG:HG3	1.89	0.54
53:D5:52:LYS:H	53:D5:53:PRO:HD2	1.73	0.54
1:CA:1373:G:H5''	7:CG:36:LYS:HZ3	1.73	0.54
25:BC:25:THR:HG21	25:BC:81:ALA:HB1	1.90	0.54
53:B5:54:GLU:HA	53:B5:57:ARG:NH1	2.20	0.54
23:BA:2186:G:H2'	23:BA:2187:G:C8	2.42	0.54
11:CK:120:ARG:HH21	11:CK:126:ARG:NH2	2.05	0.54
23:BA:2305:A:C2	28:BF:154:GLY:HA3	2.43	0.54
23:BA:2102:U:H2'	23:BA:2103:C:C6	2.42	0.54
10:CJ:38:ILE:HB	10:CJ:71:LEU:HB3	1.88	0.54
23:BA:919:G:H5'	24:BB:81:G:H1'	1.90	0.54
23:DA:529:A:H62	23:DA:2041:U:H3	1.54	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DA:276:A:H3'	23:DA:277:C:H5''	1.89	0.54
23:DA:1819:A:H4'	23:DA:1820:U:H5''	1.88	0.54
25:DC:125:ILE:HD12	25:DC:125:ILE:H	1.72	0.54
23:DA:2726:U:H5'	23:DA:2726:U:O2	2.08	0.54
32:BJ:151:HIS:HD2	32:BJ:152:PRO:O	1.90	0.54
23:BA:1161:C:O2'	40:BR:8:GLY:HA2	2.08	0.54
42:BT:28:PHE:HE2	42:BT:92:LEU:HD11	1.71	0.54
34:BL:62:LEU:HD22	34:BL:62:LEU:H	1.73	0.54
23:DA:2393:A:H5''	34:DL:62:LEU:HD12	1.90	0.54
47:BY:2:LYS:CD	47:BY:2:LYS:N	2.69	0.54
3:CC:105:GLU:HG2	3:CC:106:VAL:N	2.17	0.54
25:BC:123:ALA:HB3	25:BC:131:LEU:HD23	1.89	0.54
25:DC:123:ALA:HB3	25:DC:131:LEU:HD23	1.88	0.54
1:CA:673:G:H5''	6:CF:87:ARG:HH12	1.72	0.54
23:DA:947:G:N2	23:DA:971:C:C2	2.76	0.54
25:BC:25:THR:HG21	25:BC:81:ALA:CB	2.38	0.54
17:CQ:40:LYS:HG2	17:CQ:41:LYS:N	2.23	0.54
13:CM:34:LEU:HD22	13:CM:39:ILE:HB	1.89	0.54
29:DG:121:ILE:HD11	29:DG:140:LYS:HB3	1.90	0.54
1:CA:950:U:H2'	1:CA:951:G:H8	1.73	0.54
23:BA:319:C:H2'	23:BA:320:A:H8	1.73	0.54
1:AA:918:A:H2'	1:AA:919:A:C8	2.42	0.54
1:AA:559:A:H4'	1:AA:560:U:H5''	1.90	0.54
4:AD:12:CYS:SG	4:AD:19:LEU:HB2	2.48	0.54
25:BC:246:PRO:HD2	25:BC:255:LYS:HD3	1.89	0.54
23:DA:2293:C:H4'	37:DO:93:LYS:NZ	2.23	0.54
20:CT:29:LYS:O	20:CT:33:ILE:HG12	2.08	0.54
1:CA:1281:U:H4'	1:CA:1282:C:OP2	2.07	0.54
44:DV:110:GLY:HA3	44:DV:174:VAL:HG11	1.90	0.54
13:CM:15:VAL:O	13:CM:19:LEU:HD23	2.07	0.54
23:BA:2726:U:H5'	23:BA:2726:U:O2	2.08	0.54
1:AA:563:A:N3	1:AA:563:A:H2'	2.23	0.54
23:BA:929:G:H8	23:BA:929:G:O5'	1.90	0.54
2:CB:22:LYS:HZ3	2:CB:22:LYS:H	1.56	0.54
23:DA:1647:G:H3'	23:DA:1647:G:OP2	2.06	0.54
35:BM:54:MET:HG2	35:BM:64:ILE:HD13	1.90	0.54
1:AA:512:U:H2'	1:AA:513:C:H6	1.73	0.54
45:BW:51:VAL:N	45:BW:62:LEU:HD12	2.22	0.54
5:AE:127:ASN:O	5:AE:131:ILE:HG12	2.08	0.54
23:BA:2729:G:H2'	23:BA:2730:C:C6	2.43	0.54
39:DQ:79:PHE:HE1	39:DQ:83:LEU:HD21	1.72	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
43:DU:37:VAL:HG21	43:DU:72:VAL:HG21	1.89	0.54
28:BF:60:LEU:HD11	28:BF:92:VAL:CG1	2.30	0.54
37:BO:35:ILE:CG1	37:BO:101:LEU:HD21	2.34	0.54
23:DA:2744:G:N2	29:DG:143:GLN:HE22	1.99	0.54
1:AA:80:G:H2'	1:AA:81:G:C8	2.43	0.54
1:CA:80:G:H2'	1:CA:81:G:C8	2.43	0.54
2:AB:51:LEU:O	2:AB:55:PHE:HD2	1.91	0.54
30:BH:101:LEU:HG	30:BH:107:ILE:HG23	1.90	0.54
1:AA:405:U:H3'	1:AA:406:G:H5'	1.89	0.54
6:CF:61:LEU:HB3	6:CF:63:TYR:HE2	1.72	0.54
41:BS:73:ALA:O	41:BS:106:ILE:HG12	2.07	0.54
23:BA:2731:G:C6	23:BA:2732:G:O6	2.61	0.54
15:CO:56:LEU:O	15:CO:60:VAL:HG23	2.07	0.54
16:CP:49:LEU:HD12	16:CP:50:LYS:H	1.73	0.54
24:BB:49:C:OP1	37:BO:97:ARG:HG3	2.08	0.54
3:AC:50:ALA:HB2	3:AC:75:VAL:HB	1.89	0.54
1:CA:116:A:H61	1:CA:313:A:H1'	1.72	0.54
48:DZ:17:LYS:HD3	48:DZ:17:LYS:C	2.28	0.54
1:AA:1251:A:H2'	1:AA:1252:A:C8	2.43	0.54
23:DA:2422:A:N7	53:D5:31:HIS:CE1	2.76	0.54
42:BT:39:ILE:O	42:BT:43:VAL:HG12	2.08	0.54
49:D1:59:VAL:HG12	49:D1:60:GLU:N	2.19	0.54
25:BC:131:LEU:HA	25:BC:190:TYR:CE2	2.43	0.54
27:BE:53:THR:HG23	27:BE:55:GLY:N	2.18	0.54
26:BD:25:VAL:HG12	26:BD:181:LEU:HD12	1.90	0.54
30:DH:77:LEU:HD11	30:DH:101:LEU:HB2	1.90	0.54
8:CH:86:ILE:CB	8:CH:133:LEU:HD22	2.38	0.54
1:AA:1369:C:H2'	1:AA:1370:G:H8	1.72	0.54
33:DK:2:ILE:HG12	33:DK:8:LEU:HD11	1.88	0.54
43:DU:76:CYS:SG	43:DU:77:PRO:HD2	2.48	0.54
29:DG:109:PHE:CE1	29:DG:152:ARG:HD3	2.43	0.54
23:DA:605:C:H1'	23:DA:657:U:O2'	2.08	0.54
20:AT:29:LYS:O	20:AT:33:ILE:HG12	2.07	0.54
23:DA:137(B):G:H2'	23:DA:139:G:N7	2.22	0.54
36:BN:96:ARG:HH22	36:BN:117:VAL:HG23	1.73	0.54
23:DA:614:U:H4'	23:DA:615:G:H5''	1.90	0.54
8:CH:58:TYR:O	8:CH:59:LEU:HD23	2.08	0.54
1:AA:323:U:H4'	20:AT:22:ARG:HB3	1.90	0.54
24:BB:40:U:O5'	24:BB:40:U:H6	1.91	0.54
39:DQ:92:ARG:CD	39:DQ:94:ASN:HB3	2.38	0.53
39:BQ:92:ARG:CD	39:BQ:94:ASN:HB3	2.38	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BA:955:C:OP2	35:BM:14:ARG:HD2	2.08	0.53
17:CQ:86:GLU:O	17:CQ:90:ILE:HG12	2.07	0.53
25:BC:132:PRO:HD3	25:BC:190:TYR:CZ	2.44	0.53
25:DC:25:THR:HG21	25:DC:81:ALA:CB	2.37	0.53
25:DC:71:ASP:CB	25:DC:103:ARG:HH22	2.20	0.53
2:CB:80:ILE:HD11	2:CB:208:ILE:HG23	1.88	0.53
34:BL:18:ARG:HB3	34:BL:18:ARG:NH1	2.22	0.53
35:BM:45:GLN:H	35:BM:45:GLN:CD	2.10	0.53
37:DO:87:PHE:CE2	37:DO:89:ARG:HA	2.43	0.53
8:AH:89:PRO:HA	8:AH:92:ARG:HH11	1.72	0.53
35:BM:112:GLU:H	35:BM:112:GLU:CD	2.11	0.53
9:AI:17:VAL:HG13	9:AI:63:ILE:HD11	1.90	0.53
17:AQ:40:LYS:HG2	17:AQ:41:LYS:N	2.23	0.53
2:CB:158:LEU:HD12	2:CB:158:LEU:N	2.23	0.53
28:BF:47:LYS:HG3	28:BF:82:LEU:CD2	2.39	0.53
28:BF:134:GLY:C	28:BF:135:LEU:HD12	2.29	0.53
34:BL:132:LYS:HD2	34:BL:132:LYS:N	2.22	0.53
13:AM:4:ILE:HA	13:AM:57:ARG:HG3	1.88	0.53
27:BE:183:VAL:O	27:BE:187:VAL:HG23	2.08	0.53
24:DB:40:U:H6	24:DB:40:U:O5'	1.92	0.53
1:CA:563:A:N3	1:CA:563:A:H2'	2.23	0.53
44:BV:110:GLY:HA3	44:BV:174:VAL:HG11	1.90	0.53
33:BK:96:THR:O	33:BK:97:ARG:C	2.47	0.53
1:AA:269:C:H2'	1:AA:270:A:C8	2.42	0.53
1:AA:137:C:O4'	16:AP:63:GLY:HA3	2.08	0.53
40:DR:22:VAL:HG12	40:DR:23:GLU:H	1.73	0.53
25:DC:233:HIS:HE1	25:DC:247:ALA:H	1.56	0.53
29:BG:27:LYS:HG2	29:BG:32:GLU:HB2	1.90	0.53
19:AS:16:LEU:O	19:AS:20:LEU:HG	2.09	0.53
23:BA:911:A:C6	35:BM:9:TYR:HE1	2.25	0.53
17:CQ:56:VAL:HG23	17:CQ:81:ARG:HG3	1.89	0.53
34:DL:58:THR:C	34:DL:60:MET:H	2.12	0.53
52:D4:8:ASN:ND2	52:D4:11:LYS:N	2.54	0.53
16:AP:20:VAL:HG23	16:AP:35:LYS:HA	1.90	0.53
37:DO:34:HIS:ND1	37:DO:54:LEU:HB2	2.22	0.53
30:DH:76:THR:HA	30:DH:141:LYS:HB2	1.89	0.53
23:BA:674:G:C1'	27:BE:74:ARG:HD3	2.37	0.53
47:DY:46:GLN:HA	47:DY:46:GLN:OE1	2.09	0.53
26:DD:47:VAL:HG21	26:DD:86:PRO:HD3	1.90	0.53
1:CA:1063:C:H3'	1:CA:1064:G:H2'	1.91	0.53
27:DE:192:LEU:HD21	27:DE:194:MET:HE3	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1157:A:H4'	1:CA:1158:C:O5'	2.08	0.53
33:BK:24:VAL:CG2	33:BK:33:ALA:HB2	2.38	0.53
1:CA:363:A:H8	12:CL:32:ARG:HH21	1.56	0.53
23:BA:1154:G:O5'	23:BA:1154:G:H8	1.90	0.53
23:BA:320:A:H2'	27:BE:136:THR:HG21	1.91	0.53
1:CA:512:U:H2'	1:CA:513:C:H6	1.73	0.53
27:DE:36:VAL:O	27:DE:40:GLN:HG3	2.07	0.53
23:BA:2032:G:O2'	26:BD:145:LYS:HE2	2.09	0.53
24:BB:37:C:H2'	37:BO:95:HIS:HE1	1.71	0.53
23:DA:17:G:H4'	39:DQ:25:TRP:CH2	2.42	0.53
1:AA:1191:A:H5''	3:AC:4:LYS:NZ	2.23	0.53
34:DL:132:LYS:HD2	34:DL:132:LYS:N	2.21	0.53
40:BR:22:VAL:HG12	40:BR:23:GLU:H	1.73	0.53
15:CO:36:ILE:HG22	15:CO:37:ASN:HD22	1.73	0.53
34:DL:84:ASN:HB3	34:DL:86:LYS:HG2	1.90	0.53
53:D5:34:TRP:CG	53:D5:35:GLN:N	2.76	0.53
34:BL:88:LEU:HD22	34:BL:114:ILE:HG21	1.90	0.53
3:AC:11:ARG:HB3	3:AC:15:THR:HB	1.89	0.53
26:DD:25:VAL:HG12	26:DD:181:LEU:HD12	1.91	0.53
30:DH:130:TYR:CD2	30:DH:132:PRO:HG3	2.43	0.53
4:AD:189:PRO:CB	4:AD:194:LEU:HD21	2.38	0.53
23:DA:2186:G:H2'	23:DA:2187:G:C8	2.42	0.53
1:AA:501:C:H2'	1:AA:502:G:C8	2.43	0.53
23:DA:581:C:H2'	23:DA:582:G:H8	1.74	0.53
23:BA:2293:C:H4'	37:BO:93:LYS:NZ	2.23	0.53
1:AA:193:C:H2'	1:AA:194:C:C6	2.42	0.53
1:CA:1251:A:H2'	1:CA:1252:A:C8	2.43	0.53
23:DA:1655:A:H1'	26:DD:113:PHE:CD2	2.43	0.53
23:DA:2814:C:O2'	50:D2:29:ILE:HG13	2.08	0.53
17:AQ:97:SER:O	17:AQ:98:LEU:HD23	2.09	0.53
3:CC:173:VAL:N	3:CC:174:PRO:HD3	2.23	0.53
23:BA:2050:C:H1'	26:BD:156:MET:HE1	1.90	0.53
1:AA:434:U:H2'	1:AA:435:C:C6	2.42	0.53
17:AQ:59:ILE:HD12	17:AQ:59:ILE:N	2.23	0.53
42:DT:23:GLU:HG3	42:DT:24:GLY:H	1.73	0.53
42:BT:31:HIS:ND1	42:BT:32:PRO:HD2	2.22	0.53
1:AA:1223:C:P	19:AS:78:ARG:HH21	2.31	0.53
43:BU:11:ASP:O	43:BU:26:LYS:HA	2.09	0.53
23:BA:2543:G:H2'	23:BA:2544:G:C8	2.44	0.53
23:DA:2481:G:O2'	23:DA:2482:G:P	2.65	0.53
30:BH:142:VAL:HG12	30:BH:143:SER:N	2.23	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1427:U:H2'	1:AA:1428:A:H8	1.73	0.53
8:CH:89:PRO:HA	8:CH:92:ARG:HH11	1.73	0.53
23:BA:1614:A:H62	41:BS:93:ALA:HB2	1.73	0.53
23:DA:2401:U:C2'	23:DA:2402:C:H5''	2.39	0.53
2:AB:178:ARG:HH21	8:AH:74:PRO:HG3	1.73	0.53
6:AF:67:MET:HB2	6:AF:68:PRO:HD2	1.90	0.53
23:BA:1692:U:O2'	23:BA:1693:U:H2'	2.08	0.53
23:BA:1809:A:H2'	23:BA:1810:A:C8	2.43	0.53
26:DD:111:ARG:HD2	26:DD:160:TYR:CE1	2.44	0.53
36:DN:104:ARG:NH1	36:DN:109:ALA:HB3	2.24	0.53
4:CD:155:LEU:O	4:CD:159:ARG:HG2	2.08	0.53
23:BA:2814:C:O2'	50:B2:29:ILE:HG13	2.08	0.53
5:AE:10:MET:HA	5:AE:32:VAL:HA	1.91	0.53
20:CT:97:ALA:O	20:CT:99:LEU:N	2.41	0.53
1:AA:1236:A:H2'	1:AA:1237:C:C6	2.44	0.53
28:DF:96:ARG:O	28:DF:99:MET:HB3	2.09	0.53
23:BA:17:G:H4'	39:BQ:25:TRP:CH2	2.43	0.53
35:BM:78:PRO:O	35:BM:79:LEU:HB2	2.09	0.53
3:CC:50:ALA:HB2	3:CC:75:VAL:HB	1.90	0.53
34:BL:62:LEU:CD2	53:B5:25:MET:HB2	2.38	0.53
32:BJ:42:GLU:HA	32:BJ:82:LYS:CB	2.32	0.53
17:CQ:82:MET:O	17:CQ:86:GLU:HG2	2.08	0.53
28:BF:41:GLN:HG2	28:BF:155:MET:CB	2.37	0.53
3:AC:182:ILE:HG12	3:AC:203:PHE:HA	1.91	0.53
53:B5:50:LEU:HB2	53:B5:54:GLU:HG3	1.91	0.53
16:AP:22:THR:HA	16:AP:33:ILE:HG12	1.90	0.53
47:BY:16:LEU:HB2	47:BY:20:GLU:CG	2.35	0.53
25:BC:244:ARG:HG3	25:BC:245:PRO:N	2.24	0.53
2:CB:205:ASP:O	2:CB:211:ILE:HD11	2.09	0.53
28:DF:32:PRO:HA	28:DF:162:THR:OG1	2.09	0.53
37:BO:87:PHE:CE2	37:BO:89:ARG:HA	2.43	0.53
10:CJ:13:HIS:HB3	10:CJ:68:HIS:CD2	2.44	0.53
4:AD:104:VAL:HG11	4:AD:146:ILE:CD1	2.38	0.53
8:AH:88:LYS:HB3	8:AH:89:PRO:HD2	1.91	0.53
4:CD:104:VAL:HG11	4:CD:146:ILE:CD1	2.38	0.53
1:CA:413:G:H21	1:CA:428:G:H1'	1.73	0.53
38:BP:57:PHE:O	38:BP:59:THR:N	2.41	0.53
23:DA:814:C:O2'	23:DA:815:C:H5'	2.08	0.53
24:DB:49:C:OP1	37:DO:97:ARG:HG3	2.08	0.53
23:BA:176:G:O2'	23:BA:177:G:H5'	2.09	0.53
1:AA:1281:U:H4'	1:AA:1282:C:OP2	2.07	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DA:1932:A:H3'	23:DA:1933:G:H8	1.73	0.53
23:DA:1266:G:H5''	50:D2:23:HIS:NE2	2.23	0.53
1:CA:20:U:H2'	1:CA:21:G:O4'	2.09	0.53
1:CA:1236:A:H2'	1:CA:1237:C:C6	2.43	0.53
36:DN:96:ARG:HH22	36:DN:117:VAL:HG23	1.73	0.53
23:DA:483:A:H4'	43:DU:49:VAL:HG23	1.89	0.53
43:DU:59:GLY:C	43:DU:61:ILE:H	2.11	0.53
14:CN:48:ALA:HB2	14:CN:53:LEU:HD12	1.90	0.53
39:BQ:92:ARG:NH2	40:BR:11:GLN:H	2.06	0.53
34:DL:62:LEU:HD22	34:DL:62:LEU:H	1.74	0.53
1:AA:939:G:H2'	1:AA:940:C:C6	2.44	0.53
38:BP:26:ASP:HB2	38:BP:90:GLN:O	2.09	0.53
53:D5:50:LEU:HB2	53:D5:54:GLU:HG3	1.90	0.53
1:AA:78:G:H2'	1:AA:79:G:C8	2.44	0.53
5:AE:33:VAL:HG11	5:AE:109:ILE:HD13	1.89	0.53
6:AF:16:GLN:HA	6:AF:19:LEU:HB3	1.89	0.53
44:DV:108:PRO:HA	44:DV:142:SER:O	2.07	0.53
14:CN:24:CYS:O	14:CN:28:GLY:HA2	2.09	0.53
3:AC:35:GLU:HA	3:AC:38:ARG:HG2	1.91	0.53
4:CD:122:ARG:O	4:CD:122:ARG:HD3	2.08	0.53
39:DQ:58:ARG:O	39:DQ:62:ILE:HG12	2.08	0.53
1:AA:1152:A:OP1	10:AJ:68:HIS:CD2	2.62	0.53
23:DA:1309:G:H3'	52:D4:9:ARG:NH1	2.22	0.53
2:CB:178:ARG:HH21	8:CH:74:PRO:HG3	1.72	0.53
23:BA:1309:G:H3'	52:B4:9:ARG:NH1	2.23	0.53
23:BA:605:C:H1'	23:BA:657:U:O2'	2.09	0.53
34:DL:27:HIS:HE1	40:DR:83:ARG:HH12	1.56	0.53
6:CF:5:GLU:HG3	6:CF:93:SER:OG	2.09	0.53
23:DA:1692:U:O2'	23:DA:1693:U:H2'	2.09	0.53
34:BL:52:GLU:OE1	34:BL:52:GLU:HA	2.08	0.53
23:DA:1862:G:H2'	23:DA:1863:G:H8	1.73	0.53
20:CT:81:LYS:O	20:CT:85:MET:HG2	2.09	0.53
25:BC:242:ARG:H	25:BC:242:ARG:CD	2.07	0.53
32:DJ:157:ARG:N	32:DJ:158:PRO:CD	2.65	0.53
28:BF:107:LEU:HA	28:BF:111:LEU:HD12	1.91	0.53
33:BK:68:GLU:HB3	33:BK:78:ARG:HB2	1.89	0.53
26:BD:9:VAL:HG13	26:BD:25:VAL:O	2.09	0.53
14:AN:24:CYS:HB3	14:AN:29:ARG:N	2.21	0.53
44:BV:108:PRO:HA	44:BV:142:SER:O	2.08	0.53
26:DD:33:VAL:HG23	26:DD:47:VAL:HG13	1.91	0.53
29:BG:109:PHE:CE1	29:BG:152:ARG:HD3	2.44	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
24:BB:30:C:H2'	24:BB:31:C:H5'	1.90	0.53
1:CA:626:U:H2'	1:CA:627:G:C8	2.44	0.53
22:CV:6168:G:H2'	22:CV:6169:U:H6	1.74	0.53
38:BP:1:MET:O	38:BP:3:ARG:N	2.41	0.53
25:DC:72:LYS:HE3	25:DC:101:GLU:HG2	1.91	0.53
32:BJ:49:LEU:O	32:BJ:53:ILE:HG13	2.09	0.53
41:DS:103:ILE:H	41:DS:103:ILE:HD12	1.74	0.53
41:BS:80:PRO:O	41:BS:100:THR:HG22	2.09	0.53
35:BM:48:GLU:O	35:BM:52:VAL:HG12	2.08	0.53
23:DA:1386:C:H2'	23:DA:1387:C:C6	2.44	0.53
23:DA:1386:C:H2'	23:DA:1387:C:H6	1.74	0.53
1:CA:1125:U:H6	1:CA:1125:U:O5'	1.92	0.53
1:CA:1317:C:C2	14:CN:16:PHE:CZ	2.96	0.53
5:AE:72:GLN:O	5:AE:75:THR:HG22	2.09	0.53
23:DA:2394:C:H2'	23:DA:2395:C:C6	2.44	0.53
23:BA:637:A:OP1	34:BL:133:SER:HB3	2.09	0.53
17:AQ:80:GLY:O	17:AQ:81:ARG:HG2	2.09	0.53
29:DG:149:ARG:HA	29:DG:162:ILE:HG12	1.90	0.53
44:DV:76:LEU:H	44:DV:76:LEU:HD12	1.73	0.53
1:AA:942:G:H21	9:AI:124:GLN:NE2	2.00	0.53
11:CK:43:SER:HA	11:CK:47:VAL:HG21	1.91	0.53
25:DC:131:LEU:HA	25:DC:190:TYR:CE2	2.43	0.53
1:CA:939:G:H2'	1:CA:940:C:C6	2.44	0.53
53:B5:53:PRO:HB2	53:B5:57:ARG:HH21	1.72	0.53
43:DU:11:ASP:O	43:DU:26:LYS:HA	2.09	0.53
23:DA:2543:G:H2'	23:DA:2544:G:C8	2.44	0.53
23:DA:322:A:H3'	27:DE:169:ASN:HD21	1.72	0.53
7:AG:15:ASP:HA	7:AG:24:THR:HG23	1.91	0.53
26:BD:47:VAL:HG21	26:BD:86:PRO:HD3	1.90	0.53
23:DA:597:U:O2'	34:DL:15:ARG:HG2	2.09	0.53
8:CH:97:VAL:HG13	8:CH:98:LYS:N	2.24	0.53
23:BA:1257:C:H4'	27:BE:83:PHE:CE2	2.44	0.53
23:DA:2648:C:H2'	23:DA:2649:U:C6	2.43	0.53
1:CA:238:G:P	17:CQ:25:ARG:HH22	2.32	0.53
1:CA:1154:G:H2'	1:CA:1155:G:H8	1.74	0.53
39:BQ:8:VAL:HG11	39:BQ:12:ARG:CZ	2.39	0.53
46:DX:45:ASN:ND2	46:DX:47:GLN:HE21	2.07	0.53
1:AA:1317:C:C2	14:AN:16:PHE:CZ	2.97	0.53
1:AA:765:G:H5''	1:AA:766:A:OP1	2.09	0.53
38:DP:1:MET:O	38:DP:3:ARG:N	2.40	0.53
20:AT:81:LYS:O	20:AT:85:MET:HG2	2.09	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:963:G:H2'	1:AA:964:A:C8	2.44	0.53
23:BA:1862:G:H2'	23:BA:1863:G:H8	1.74	0.53
1:AA:1125:U:O5'	1:AA:1125:U:H6	1.92	0.53
23:DA:127:A:H5''	23:DA:128:C:O4'	2.09	0.53
17:CQ:97:SER:O	17:CQ:98:LEU:HD23	2.09	0.53
1:CA:149:A:H2'	1:CA:150:C:C6	2.43	0.53
23:BA:2056:G:N2	23:BA:2057:A:C1'	2.72	0.53
23:BA:1971:A:C5	25:BC:241:PRO:HG3	2.44	0.53
53:B5:34:TRP:CG	53:B5:35:GLN:N	2.76	0.53
49:D1:42:CYS:SG	49:D1:46:ASN:HB3	2.49	0.53
35:BM:140:ALA:HB3	44:BV:53:ILE:HD13	1.90	0.53
35:BM:141:GLN:HA	44:BV:71:VAL:O	2.08	0.53
25:DC:132:PRO:HD3	25:DC:190:TYR:CZ	2.44	0.53
9:CI:65:VAL:HG21	9:CI:73:GLN:HB3	1.91	0.53
23:DA:2822:G:O6	36:DN:4:LEU:HD23	2.08	0.53
12:CL:82:VAL:HG22	12:CL:83:LEU:N	2.24	0.53
7:CG:69:VAL:CA	7:CG:138:LYS:HD2	2.35	0.53
34:DL:17:LYS:O	34:DL:19:VAL:HG22	2.09	0.53
46:BX:27:GLU:CB	46:BX:33:LYS:HG3	2.39	0.53
24:BB:104:A:O4'	44:BV:29:TYR:HE1	1.92	0.53
1:CA:1356:G:H2'	1:CA:1357:A:H8	1.74	0.53
23:DA:1614:A:H61	41:DS:88:ARG:H	1.55	0.53
10:CJ:54:PHE:HD2	10:CJ:55:LYS:HG3	1.73	0.53
23:BA:2401:U:C2'	23:BA:2402:C:H5''	2.38	0.53
6:AF:12:PRO:HG2	6:AF:55:ASP:OD2	2.09	0.53
1:CA:59:A:H3'	1:CA:331:G:H22	1.74	0.53
22:CV:6168:G:H2'	22:CV:6169:U:C6	2.44	0.53
37:DO:41:ASP:OD2	37:DO:44:LYS:HD3	2.09	0.53
23:DA:2731:G:C6	23:DA:2732:G:O6	2.62	0.53
44:DV:118:GLN:HB2	44:DV:173:ALA:O	2.09	0.53
1:AA:950:U:H2'	1:AA:951:G:H8	1.74	0.53
23:BA:1028:A:N6	23:BA:1125:G:H2'	2.24	0.53
34:BL:75:ILE:HD13	34:BL:77:ARG:NE	2.24	0.53
35:BM:68:ILE:HG23	35:BM:103:MET:HA	1.91	0.53
41:DS:110:LYS:HG3	41:DS:111:HIS:ND1	2.24	0.53
6:AF:5:GLU:HB3	6:AF:62:TRP:HE1	1.73	0.53
1:AA:971:G:H1'	1:AA:1365:G:O2'	2.08	0.53
40:DR:38:LEU:HD22	40:DR:52:VAL:HG11	1.91	0.53
23:DA:1164:G:C5'	23:DA:1164:G:H8	2.22	0.53
23:BA:2422:A:N7	53:B5:31:HIS:CE1	2.77	0.53
23:DA:637:A:OP1	34:DL:133:SER:HB3	2.09	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:735:C:H2'	1:CA:736:C:C6	2.44	0.53
1:AA:735:C:H2'	1:AA:736:C:C6	2.42	0.53
49:B1:42:CYS:SG	49:B1:46:ASN:HB3	2.48	0.53
28:DF:107:LEU:HA	28:DF:111:LEU:HD12	1.91	0.53
9:CI:10:ARG:HH21	9:CI:107:ARG:HB2	1.74	0.53
23:DA:948:G:H8	23:DA:948:G:C5'	2.18	0.53
23:BA:2481:G:O2'	23:BA:2482:G:P	2.67	0.53
23:DA:323:G:H5'	27:DE:169:ASN:HD21	1.73	0.53
9:AI:10:ARG:HH21	9:AI:107:ARG:HB2	1.74	0.53
12:AL:65:VAL:HG11	12:AL:97:TYR:CE1	2.44	0.53
36:BN:12:ARG:HD3	36:BN:16:HIS:ND1	2.24	0.53
6:CF:67:MET:HB2	6:CF:68:PRO:HD2	1.90	0.53
1:AA:161:A:H2'	1:AA:162:A:C8	2.44	0.53
12:CL:32:ARG:HA	12:CL:32:ARG:HE	1.71	0.53
23:BA:1386:C:H2'	23:BA:1387:C:H6	1.74	0.53
23:BA:55:G:H2'	23:BA:56:A:H8	1.73	0.53
9:AI:83:ARG:HA	9:AI:86:VAL:HG12	1.91	0.53
23:DA:2010:G:H5''	41:DS:42:ARG:HB2	1.91	0.53
2:CB:17:PHE:HB2	2:CB:42:ILE:CG2	2.40	0.53
30:DH:117:GLU:HG3	30:DH:118:LYS:N	2.24	0.53
27:BE:18:ARG:HG3	27:BE:18:ARG:O	2.08	0.53
23:DA:486:C:H4'	41:DS:60:ASN:HD22	1.74	0.53
13:CM:87:TYR:O	13:CM:91:ARG:HG2	2.08	0.53
2:AB:17:PHE:HB2	2:AB:42:ILE:CG2	2.39	0.53
40:BR:6:LYS:O	40:BR:37:VAL:HG21	2.10	0.52
15:AO:36:ILE:HD12	15:AO:63:ARG:HH11	1.74	0.52
19:CS:16:LEU:O	19:CS:20:LEU:HG	2.09	0.52
42:DT:39:ILE:O	42:DT:43:VAL:HG12	2.09	0.52
27:BE:63:LYS:HE3	27:BE:75:HIS:O	2.09	0.52
23:BA:270(L):C:H2'	23:BA:270(M):U:H5''	1.91	0.52
43:BU:95:LYS:HG2	43:BU:100:ALA:HA	1.91	0.52
37:BO:35:ILE:O	37:BO:53:SER:HB2	2.09	0.52
23:DA:1510:A:H2'	23:DA:1511:A:H8	1.71	0.52
36:DN:10:LEU:HD22	36:DN:17:ARG:CD	2.39	0.52
7:CG:15:ASP:HA	7:CG:24:THR:HG23	1.91	0.52
34:BL:143:GLY:C	34:BL:145:PRO:HD3	2.30	0.52
1:AA:501:C:H2'	1:AA:502:G:H8	1.72	0.52
45:DW:48:GLY:HA3	45:DW:80:HIS:ND1	2.24	0.52
3:AC:173:VAL:N	3:AC:174:PRO:HD3	2.24	0.52
1:CA:160:A:H2'	1:CA:161:A:O4'	2.08	0.52
1:CA:865:A:H5'	1:CA:1078:U:O4	2.08	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
13:AM:10:PRO:HB2	13:AM:18:ALA:HB1	1.90	0.52
17:AQ:55:ASP:HB3	17:AQ:76:LEU:HD13	1.92	0.52
3:AC:184:TYR:HE2	3:AC:186:PHE:HB2	1.74	0.52
13:AM:17:VAL:HG12	13:AM:21:TYR:HE1	1.74	0.52
23:DA:263:C:H2'	23:DA:264:C:O4'	2.09	0.52
23:DA:176:G:O2'	23:DA:177:G:H5'	2.09	0.52
1:CA:772:U:H2'	1:CA:773:G:O4'	2.09	0.52
1:AA:1053:G:N7	1:AA:1200:C:H5''	2.24	0.52
4:CD:142:PRO:HA	4:CD:185:PHE:HD2	1.74	0.52
37:BO:41:ASP:OD2	37:BO:44:LYS:HD3	2.09	0.52
1:CA:1229:A:H2'	1:CA:1230:C:C6	2.44	0.52
13:AM:87:TYR:O	13:AM:91:ARG:HG2	2.09	0.52
23:BA:676:A:H2	23:BA:802:A:H61	1.57	0.52
17:CQ:80:GLY:O	17:CQ:81:ARG:HG2	2.09	0.52
23:BA:2210:G:H21	23:BA:2211:G:C5'	2.23	0.52
30:DH:77:LEU:HG	30:DH:101:LEU:HD13	1.91	0.52
8:CH:88:LYS:HB3	8:CH:89:PRO:HD2	1.90	0.52
23:DA:2358:G:C6	23:DA:2359:C:C4	2.97	0.52
23:DA:2271:G:H2'	23:DA:2272:U:C6	2.44	0.52
20:AT:26:ASN:HD22	20:AT:27:LYS:N	2.07	0.52
33:BK:2:ILE:HG12	33:BK:8:LEU:HD11	1.91	0.52
1:CA:601:C:H2'	1:CA:602:A:H8	1.73	0.52
2:AB:158:LEU:HD12	2:AB:158:LEU:N	2.24	0.52
28:DF:47:LYS:HG3	28:DF:82:LEU:CD2	2.38	0.52
1:CA:17:U:H2'	1:CA:18:C:C6	2.45	0.52
45:BW:48:GLY:HA3	45:BW:80:HIS:ND1	2.24	0.52
10:AJ:30:SER:HB2	10:AJ:80:LYS:CG	2.40	0.52
5:CE:10:MET:HA	5:CE:32:VAL:HA	1.91	0.52
1:AA:1154:G:H2'	1:AA:1155:G:H8	1.74	0.52
23:BA:263:C:H2'	23:BA:264:C:O4'	2.09	0.52
23:BA:498:G:N3	43:BU:47:LYS:HE3	2.24	0.52
14:AN:48:ALA:HB2	14:AN:53:LEU:HD12	1.91	0.52
8:AH:80:ILE:N	8:AH:80:ILE:HD12	2.25	0.52
23:DA:593:G:O2'	53:D5:62:LEU:HD13	2.09	0.52
39:DQ:88:ILE:O	39:DQ:88:ILE:HG13	2.09	0.52
23:DA:1022:G:H8	32:DJ:92:GLN:HE22	1.56	0.52
47:BY:2:LYS:HA	47:BY:5:GLU:OE2	2.08	0.52
11:CK:29:ILE:C	11:CK:29:ILE:HD12	2.30	0.52
16:AP:21:VAL:HG23	16:AP:33:ILE:HB	1.91	0.52
25:BC:71:ASP:HB3	25:BC:103:ARG:NH2	2.20	0.52
12:AL:82:VAL:HG22	12:AL:83:LEU:N	2.24	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DA:2517:C:C6	23:DA:2542:A:C2	2.96	0.52
13:CM:3:ARG:HH21	13:CM:7:VAL:HG13	1.74	0.52
36:BN:10:LEU:HD22	36:BN:17:ARG:CD	2.40	0.52
36:BN:21:TYR:HE2	36:BN:43:GLU:HB3	1.73	0.52
2:CB:51:LEU:O	2:CB:55:PHE:HD2	1.91	0.52
30:BH:76:THR:HA	30:BH:141:LYS:HB2	1.90	0.52
23:DA:775:G:C4	23:DA:794:G:C8	2.97	0.52
35:BM:60:ARG:H	44:BV:179:ASP:CB	2.22	0.52
1:CA:1369:C:H2'	1:CA:1370:G:H8	1.72	0.52
1:AA:1014:A:H2	1:AA:1219:U:H1'	1.75	0.52
23:DA:568:U:O4	40:DR:78:LYS:NZ	2.40	0.52
23:BA:2358:G:C6	23:BA:2359:C:C4	2.97	0.52
10:AJ:13:HIS:HB3	10:AJ:68:HIS:CD2	2.45	0.52
9:AI:113:LYS:HG2	9:AI:119:ALA:HA	1.91	0.52
23:BA:1276:A:O2'	36:BN:16:HIS:HE1	1.92	0.52
12:CL:40:ARG:HD3	12:CL:41:THR:O	2.09	0.52
35:DM:58:PHE:O	35:DM:58:PHE:HD1	1.92	0.52
1:CA:512:U:H2'	1:CA:513:C:C6	2.44	0.52
5:AE:10:MET:HG3	5:AE:13:ILE:HD11	1.90	0.52
23:BA:118:A:N3	23:BA:178:G:H1'	2.25	0.52
4:CD:100:ARG:NH1	4:CD:137:SER:HA	2.24	0.52
35:DM:78:PRO:O	35:DM:79:LEU:HB2	2.09	0.52
30:BH:66:GLU:HB3	30:BH:67:ARG:NH1	2.24	0.52
1:CA:332:G:OP2	20:CT:10:LEU:HD23	2.10	0.52
44:DV:94:GLU:H	44:DV:94:GLU:CD	2.12	0.52
41:BS:110:LYS:HG3	41:BS:111:HIS:ND1	2.25	0.52
38:BP:6:LEU:O	38:BP:10:VAL:HG23	2.09	0.52
1:CA:765:G:H5''	1:CA:766:A:OP1	2.09	0.52
23:BA:1790:C:O2'	25:BC:209:ALA:HB2	2.09	0.52
40:BR:15:GLU:HB3	40:BR:16:PRO:HD2	1.92	0.52
37:DO:49:VAL:HG11	37:DO:73:LEU:HA	1.92	0.52
17:CQ:54:GLY:O	17:CQ:81:ARG:HB2	2.09	0.52
38:BP:48:ILE:HD12	38:BP:48:ILE:H	1.73	0.52
53:B5:50:LEU:HD13	53:B5:57:ARG:CZ	2.40	0.52
30:BH:72:LEU:HD12	30:BH:140:LEU:HD13	1.91	0.52
23:BA:1173:G:H3'	23:BA:1174:A:C5'	2.40	0.52
33:DK:68:GLU:H	33:DK:68:GLU:CD	2.12	0.52
11:AK:21:ILE:HD12	11:AK:21:ILE:N	2.25	0.52
23:DA:779:U:P	25:DC:49:ILE:HG13	2.49	0.52
1:AA:1347:G:C8	9:AI:107:ARG:HB3	2.44	0.52
9:AI:65:VAL:HG21	9:AI:73:GLN:HB3	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:BD:33:VAL:HG23	26:BD:47:VAL:HG13	1.90	0.52
4:AD:166:LYS:O	4:AD:166:LYS:HD2	2.10	0.52
43:DU:90:LEU:HG	43:DU:91:GLU:HG2	1.91	0.52
52:D4:21:ARG:HB3	52:D4:31:LEU:CD2	2.40	0.52
38:DP:57:PHE:HE2	38:DP:79:HIS:HB2	1.74	0.52
23:BA:1567:A:H2'	25:BC:84:TYR:HE2	1.75	0.52
39:DQ:49:HIS:HA	39:DQ:52:ARG:HB2	1.92	0.52
44:DV:5:LEU:HB3	44:DV:59:LEU:HD23	1.92	0.52
23:BA:1817:G:OP1	25:BC:88:ARG:NH2	2.42	0.52
1:CA:350:G:O2'	1:CA:351:G:H5'	2.08	0.52
23:BA:2774:C:H2'	23:BA:2775:A:O4'	2.09	0.52
28:BF:96:ARG:O	28:BF:99:MET:HB3	2.09	0.52
1:AA:1260:C:H4'	1:AA:1284:C:H5'	1.92	0.52
23:DA:1826:G:OP1	25:DC:233:HIS:HD2	1.92	0.52
15:CO:36:ILE:HD12	15:CO:63:ARG:HH11	1.74	0.52
23:BA:1164:G:H8	23:BA:1164:G:C5'	2.22	0.52
34:DL:62:LEU:CD2	53:D5:25:MET:HB2	2.38	0.52
47:BY:12:GLU:C	47:BY:14:ARG:H	2.13	0.52
23:DA:2439:A:H8	23:DA:2439:A:C5'	2.22	0.52
2:AB:205:ASP:O	2:AB:211:ILE:HD11	2.09	0.52
34:BL:58:THR:C	34:BL:60:MET:H	2.12	0.52
27:DE:53:THR:HG23	27:DE:55:GLY:N	2.19	0.52
27:DE:34:TRP:HB2	34:DL:10:PRO:O	2.10	0.52
27:BE:29:ASN:H	27:BE:112:MET:CE	2.22	0.52
34:DL:143:GLY:C	34:DL:145:PRO:HD3	2.29	0.52
1:CA:1504:G:OP1	1:CA:1507:A:H4'	2.10	0.52
1:CA:817:C:H1'	1:CA:819:A:H5'	1.91	0.52
23:BA:1386:C:H2'	23:BA:1387:C:C6	2.44	0.52
5:CE:10:MET:HG3	5:CE:13:ILE:HD11	1.90	0.52
6:AF:14:LEU:HD21	6:AF:18:GLN:HB2	1.92	0.52
1:CA:909:A:OP1	12:CL:20:LYS:HD2	2.08	0.52
1:AA:136(B):C:H5'	1:AA:1363:A:H2'	1.90	0.52
28:DF:139:LEU:HA	28:DF:144:ILE:HG21	1.91	0.52
23:BA:127:A:H5''	23:BA:128:C:O4'	2.09	0.52
27:BE:179:GLU:CD	27:BE:179:GLU:H	2.12	0.52
1:AA:164:U:H2'	1:AA:165:C:C6	2.45	0.52
10:CJ:30:SER:HB2	10:CJ:80:LYS:CG	2.39	0.52
33:DK:112:MET:HA	33:DK:115:VAL:HG22	1.91	0.52
1:CA:1240:U:OP1	7:CG:115:ARG:HA	2.09	0.52
25:BC:242:ARG:CD	25:BC:242:ARG:N	2.70	0.52
15:AO:36:ILE:HG22	15:AO:37:ASN:HD22	1.74	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:DL:88:LEU:HD22	34:DL:114:ILE:HG21	1.90	0.52
23:DA:959:A:O2'	23:DA:960:A:H5'	2.10	0.52
29:BG:149:ARG:HA	29:BG:162:ILE:HG12	1.91	0.52
23:BA:2712:U:H1'	23:BA:712(B):A:H8	1.66	0.52
17:AQ:81:ARG:HA	17:AQ:81:ARG:HE	1.75	0.52
38:DP:26:ASP:HB2	38:DP:90:GLN:O	2.09	0.52
3:CC:182:ILE:HG12	3:CC:203:PHE:HA	1.92	0.52
16:CP:20:VAL:HG23	16:CP:35:LYS:HA	1.90	0.52
26:BD:201:THR:CG2	26:BD:202:LYS:N	2.73	0.52
47:DY:39:ALA:HA	47:DY:45:SER:CB	2.39	0.52
14:AN:24:CYS:O	14:AN:28:GLY:HA2	2.09	0.52
35:DM:60:ARG:H	44:DV:179:ASP:CB	2.22	0.52
30:DH:6:LEU:HA	30:DH:15:VAL:HG13	1.90	0.52
1:AA:413:G:H21	1:AA:428:G:H1'	1.73	0.52
1:AA:976:G:H8	1:AA:1358:U:H2'	1.72	0.52
8:CH:31:PHE:O	8:CH:35:ILE:HG12	2.10	0.52
23:DA:871:U:H4'	35:DM:69:PHE:CE2	2.45	0.52
40:DR:15:GLU:HB3	40:DR:16:PRO:HD2	1.92	0.52
23:BA:1027:A:C2	23:BA:2488:A:H5'	2.45	0.52
23:DA:195:A:H61	23:DA:198:C:H3'	1.75	0.52
25:BC:233:HIS:HE1	25:BC:247:ALA:H	1.58	0.52
37:DO:35:ILE:O	37:DO:53:SER:HB2	2.10	0.52
49:D1:46:ASN:HB2	49:D1:64:LYS:HB2	1.90	0.52
17:AQ:54:GLY:O	17:AQ:81:ARG:HB2	2.10	0.52
23:DA:270(L):C:H2'	23:DA:270(M):U:H5''	1.91	0.52
6:AF:87:ARG:HG2	6:AF:87:ARG:HH11	1.74	0.52
25:BC:31:LYS:HE3	25:BC:33:LEU:HD21	1.92	0.52
47:BY:39:ALA:HA	47:BY:45:SER:CB	2.39	0.52
28:BF:85:GLY:C	28:BF:86:MET:HG3	2.30	0.52
23:BA:819:A:C4	23:BA:1189:A:C2	2.97	0.52
34:DL:16:ARG:C	34:DL:16:ARG:HE	2.13	0.52
8:AH:86:ILE:CB	8:AH:133:LEU:HD22	2.38	0.52
25:DC:204:ILE:HD12	25:DC:204:ILE:O	2.10	0.52
23:BA:1953:A:C2	23:BA:2549:G:N3	2.78	0.52
23:BA:814:C:O2'	23:BA:815:C:H5'	2.09	0.52
23:DA:830:G:H4'	23:DA:831:G:OP2	2.09	0.52
1:CA:715:A:H2'	1:CA:716:A:C8	2.45	0.52
29:BG:121:ILE:HD11	29:BG:140:LYS:HB3	1.90	0.52
1:CA:626:U:H2'	1:CA:627:G:H8	1.75	0.52
15:CO:39:LEU:HD12	15:CO:56:LEU:HB2	1.92	0.52
23:DA:144:C:H2'	23:DA:145:G:C8	2.45	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DA:298:G:P	43:DU:85:VAL:HG22	2.49	0.52
23:DA:553:U:O2'	23:DA:554:U:H5'	2.10	0.52
23:BA:49:A:H5''	23:BA:51:G:O4'	2.10	0.52
1:CA:721:G:H4'	1:CA:722:A:O4'	2.09	0.52
6:AF:21:LEU:O	6:AF:25:ILE:HG12	2.10	0.52
2:AB:74:LYS:O	2:AB:78:GLN:HG3	2.10	0.52
27:DE:18:ARG:HG3	27:DE:18:ARG:O	2.09	0.52
1:AA:488:C:O5'	1:AA:488:C:H6	1.93	0.52
23:DA:1592:C:H2'	23:DA:1593:G:H8	1.75	0.52
1:CA:1053:G:N7	1:CA:1200:C:H5''	2.24	0.52
39:DQ:92:ARG:NH2	40:DR:11:GLN:H	2.07	0.52
32:BJ:135:LEU:HD23	32:BJ:136:GLY:H	1.75	0.52
35:DM:81:VAL:CG1	35:DM:82:ARG:HG2	2.37	0.52
43:BU:96:ILE:HG23	43:BU:101:LYS:O	2.09	0.52
6:CF:87:ARG:HH11	6:CF:87:ARG:HG2	1.75	0.52
23:BA:2822:G:O6	36:BN:4:LEU:HD23	2.10	0.52
26:DD:101:ARG:HD3	26:DD:169:ASN:ND2	2.24	0.52
26:DD:201:THR:CG2	26:DD:202:LYS:N	2.72	0.52
43:DU:31:LEU:N	43:DU:31:LEU:HD23	2.23	0.52
24:BB:71:C:C2	24:BB:72:G:C8	2.97	0.52
23:DA:556:G:H2'	23:DA:557:U:H6	1.70	0.52
1:CA:1226:C:N4	13:CM:104:ARG:HB2	2.23	0.52
3:AC:19:GLU:HA	3:AC:54:ARG:HE	1.75	0.52
23:DA:2688:U:H3'	23:DA:2688:U:O2	2.09	0.52
34:BL:27:HIS:HE1	40:BR:83:ARG:HH12	1.56	0.52
40:DR:28:GLU:HB2	40:DR:31:ALA:CB	2.39	0.52
35:DM:58:PHE:CD1	35:DM:61:GLY:HA3	2.45	0.52
1:AA:255:G:H1'	17:AQ:16:GLN:NE2	2.25	0.52
1:AA:817:C:H1'	1:AA:819:A:H5'	1.90	0.52
41:DS:46:PHE:O	41:DS:50:VAL:HG12	2.09	0.52
39:BQ:8:VAL:HG11	39:BQ:12:ARG:NE	2.24	0.52
1:CA:1423:G:H2'	1:CA:1424:C:C6	2.44	0.52
41:DS:80:PRO:O	41:DS:100:THR:HG22	2.10	0.52
1:AA:216:G:H2'	1:AA:217:C:C6	2.45	0.52
25:DC:166:GLN:CA	25:DC:166:GLN:HE21	2.22	0.52
1:CA:836:G:C6	1:CA:851:G:C6	2.98	0.52
23:BA:871:U:H4'	35:BM:69:PHE:CE2	2.45	0.52
26:DD:172:VAL:HG13	26:DD:182:LEU:HD11	1.90	0.52
23:BA:298:G:P	43:BU:85:VAL:HG22	2.49	0.52
1:AA:20:U:H2'	1:AA:21:G:O4'	2.10	0.52
1:CA:501:C:H2'	1:CA:502:G:C8	2.45	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:DL:52:GLU:HA	34:DL:52:GLU:OE1	2.10	0.52
32:DJ:135:LEU:HD23	32:DJ:136:GLY:H	1.75	0.52
23:BA:1022:G:O2'	23:BA:1023:U:OP2	2.25	0.52
42:DT:30:VAL:HG12	42:DT:31:HIS:N	2.24	0.52
23:DA:2210:G:H21	23:DA:2211:G:C5'	2.22	0.52
23:DA:1568:G:OP2	25:DC:63:ARG:NH2	2.43	0.52
23:DA:1173:G:H3'	23:DA:1174:A:C5'	2.40	0.52
6:CF:16:GLN:HA	6:CF:19:LEU:HB3	1.92	0.52
5:AE:36:ASP:O	5:AE:37:ARG:HB2	2.10	0.52
23:DA:661:C:H2'	23:DA:662:G:C8	2.45	0.52
1:CA:1225:A:H5''	1:CA:1226:C:OP2	2.10	0.52
52:D4:34:ARG:HB3	52:D4:42:LEU:HD22	1.92	0.52
10:AJ:54:PHE:HD2	10:AJ:55:LYS:HG3	1.73	0.52
1:AA:1151:A:OP1	10:AJ:41:PRO:HA	2.10	0.52
23:DA:1980:G:H3'	23:DA:1981:A:C5'	2.40	0.52
1:CA:794:A:H4'	1:CA:1521:G:O2'	2.09	0.52
1:AA:1063:C:H3'	1:AA:1064:G:H2'	1.91	0.52
23:DA:2305:A:C2	28:DF:154:GLY:HA3	2.44	0.52
23:DA:1495:A:H2'	23:DA:1495:A:N3	2.25	0.52
37:DO:15:ARG:O	37:DO:19:LYS:HG3	2.10	0.52
35:BM:58:PHE:O	35:BM:58:PHE:HD1	1.91	0.52
1:AA:512:U:H2'	1:AA:513:C:C6	2.44	0.52
1:AA:17:U:H2'	1:AA:18:C:C6	2.44	0.52
6:AF:5:GLU:HG3	6:AF:93:SER:OG	2.10	0.52
9:CI:83:ARG:HA	9:CI:86:VAL:HG12	1.91	0.52
23:DA:2371:G:O2'	51:D3:45:LYS:HB3	2.10	0.52
1:CA:983:A:H3'	1:CA:983:A:N3	2.24	0.52
16:CP:8:ARG:HB3	16:CP:28:ARG:NH1	2.25	0.52
4:CD:166:LYS:O	4:CD:166:LYS:HD2	2.10	0.52
27:BE:36:VAL:O	27:BE:40:GLN:HG3	2.08	0.52
19:AS:12:ASP:HB2	19:AS:15:LEU:HD23	1.92	0.52
23:DA:2258:C:H4'	23:DA:2259:G:OP2	2.10	0.52
39:BQ:79:PHE:HE1	39:BQ:83:LEU:HD21	1.75	0.52
43:BU:81:LYS:CD	43:BU:97:ARG:HB3	2.39	0.52
23:BA:959:A:O2'	23:BA:960:A:H5'	2.09	0.52
23:DA:848:G:C4	23:DA:933:A:C8	2.95	0.52
52:B4:8:ASN:ND2	52:B4:11:LYS:N	2.53	0.52
53:B5:52:LYS:H	53:B5:53:PRO:HD2	1.75	0.52
16:AP:4:ILE:HD13	16:AP:66:PRO:HG3	1.92	0.52
16:CP:21:VAL:HG23	16:CP:33:ILE:HB	1.92	0.52
25:BC:204:ILE:O	25:BC:204:ILE:HD12	2.10	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:729:A:H2'	1:CA:730:G:C8	2.43	0.52
4:AD:21:LEU:HD12	4:AD:22:LYS:H	1.75	0.52
23:BA:917:A:H2'	23:BA:918:A:O4'	2.09	0.52
23:DA:588:U:H1'	27:DE:90:PHE:CD1	2.45	0.52
36:BN:81:ASP:O	36:BN:85:PRO:HG2	2.10	0.52
1:AA:794:A:H4'	1:AA:1521:G:O2'	2.10	0.52
23:DA:2305:A:H3'	23:DA:2306:C:H5''	1.92	0.52
1:AA:513:C:H2'	1:AA:514:C:C6	2.46	0.52
35:DM:52:VAL:O	35:DM:56:ARG:HB2	2.10	0.52
35:DM:54:MET:HG2	35:DM:64:ILE:HD13	1.91	0.52
23:DA:1790:C:O2'	25:DC:209:ALA:HB2	2.10	0.52
41:DS:19:LEU:HB3	50:D2:25:LEU:CD1	2.41	0.52
1:AA:1201:A:H4'	1:AA:1202:G:O5'	2.10	0.52
1:CA:1306:A:H2'	1:CA:1307:U:C6	2.45	0.52
1:CA:216:G:H2'	1:CA:217:C:C6	2.45	0.52
23:BA:1206:G:C6	23:BA:1207:C:C4	2.98	0.52
1:AA:1229:A:H2'	1:AA:1230:C:C6	2.44	0.52
3:CC:184:TYR:CE2	3:CC:186:PHE:HB2	2.45	0.52
25:BC:72:LYS:HE3	25:BC:101:GLU:HG2	1.92	0.52
23:BA:2094:G:N2	23:BA:2196:C:H1'	2.24	0.52
51:D3:30:THR:HG22	51:D3:31:PRO:HD2	1.92	0.52
23:BA:614:U:H4'	23:BA:615:G:H5''	1.92	0.52
23:BA:2371:G:O2'	51:B3:45:LYS:HB3	2.09	0.52
39:DQ:92:ARG:HD2	39:DQ:95:LEU:CG	2.38	0.51
39:BQ:88:ILE:HG13	39:BQ:88:ILE:O	2.10	0.51
30:DH:92:VAL:HG22	30:DH:120:ILE:HD12	1.92	0.51
17:CQ:81:ARG:HE	17:CQ:81:ARG:HA	1.75	0.51
11:AK:43:SER:HA	11:AK:47:VAL:HG21	1.92	0.51
34:DL:58:THR:C	34:DL:61:ARG:HE	2.13	0.51
1:AA:80:G:H8	1:AA:80:G:P	2.32	0.51
34:BL:58:THR:C	34:BL:61:ARG:HE	2.13	0.51
16:CP:22:THR:HA	16:CP:33:ILE:HG12	1.91	0.51
27:BE:182:ASN:O	27:BE:186:ILE:HG12	2.10	0.51
23:DA:2415:G:H4'	34:DL:66:GLY:HA2	1.92	0.51
30:DH:142:VAL:HG12	30:DH:143:SER:H	1.74	0.51
3:CC:19:GLU:HA	3:CC:54:ARG:HE	1.75	0.51
23:BA:518:G:H4'	41:BS:18:ARG:NH1	2.24	0.51
43:BU:90:LEU:HG	43:BU:91:GLU:HG2	1.92	0.51
40:BR:28:GLU:HB2	40:BR:31:ALA:CB	2.40	0.51
23:BA:2747:G:O6	23:BA:2755:C:H5''	2.10	0.51
23:DA:2747:G:O6	23:DA:2755:C:H5''	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BA:380:U:O2	46:BX:20:ARG:NH2	2.43	0.51
1:CA:513:C:H2'	1:CA:514:C:C6	2.44	0.51
23:DA:2729:G:H2'	23:DA:2730:C:C6	2.45	0.51
23:BA:1537:C:H2'	23:BA:1538:G:O4'	2.10	0.51
41:DS:84:ARG:HB2	41:DS:96:ILE:HG22	1.92	0.51
28:BF:139:LEU:HA	28:BF:144:ILE:HG21	1.91	0.51
23:DA:2094:G:N2	23:DA:2196:C:H1'	2.25	0.51
23:BA:486:C:H4'	41:BS:60:ASN:HD22	1.74	0.51
6:AF:36:ARG:HH21	6:AF:38:GLU:HG2	1.76	0.51
1:CA:663:A:H5''	18:CR:61:LYS:HE2	1.92	0.51
33:DK:24:VAL:HG23	33:DK:33:ALA:HB2	1.92	0.51
27:DE:82:ILE:HG13	27:DE:82:ILE:O	2.09	0.51
23:DA:2320:A:N3	23:DA:2320:A:H2'	2.25	0.51
23:BA:2090:G:H21	46:BX:45:ASN:ND2	2.08	0.51
23:BA:2188:C:H2'	23:BA:2189:U:O4'	2.10	0.51
23:DA:2056:G:N2	23:DA:2057:A:H1'	2.25	0.51
7:AG:113:GLU:HB3	7:AG:118:VAL:HG23	1.92	0.51
42:BT:11:PRO:HG3	47:BY:37:PHE:CE2	2.46	0.51
32:BJ:157:ARG:N	32:BJ:158:PRO:CD	2.65	0.51
2:AB:163:PHE:HD1	2:AB:185:ILE:HG13	1.75	0.51
35:BM:75:THR:HA	35:BM:88:GLY:HA3	1.89	0.51
46:DX:13:ILE:HG23	46:DX:14:VAL:H	1.75	0.51
1:CA:1267:C:O2	1:CA:1327:C:H4'	2.10	0.51
36:DN:10:LEU:HB2	36:DN:17:ARG:CZ	2.40	0.51
28:BF:74:LYS:HE3	28:BF:74:LYS:HA	1.93	0.51
33:BK:68:GLU:CD	33:BK:68:GLU:H	2.13	0.51
30:BH:107:ILE:HG13	30:BH:109:ILE:HG23	1.92	0.51
18:AR:50:ILE:HD11	18:AR:74:ARG:NH1	2.24	0.51
44:DV:104:PHE:HB3	44:DV:141:VAL:HG11	1.93	0.51
1:CA:1194:U:H2'	1:CA:1195:C:H6	1.75	0.51
23:BA:2688:U:H3'	23:BA:2688:U:O2	2.10	0.51
33:DK:2:ILE:CG1	33:DK:8:LEU:HD11	2.40	0.51
1:AA:1520:G:H2'	1:AA:1521:G:H8	1.75	0.51
38:BP:57:PHE:HE2	38:BP:79:HIS:HB2	1.76	0.51
23:DA:2105:C:H2'	23:DA:2106:G:C8	2.45	0.51
3:CC:184:TYR:HE2	3:CC:186:PHE:HB2	1.75	0.51
23:BA:2094:G:P	30:BH:22:LYS:HD2	2.50	0.51
6:CF:14:LEU:HD21	6:CF:18:GLN:HB2	1.91	0.51
44:BV:94:GLU:H	44:BV:94:GLU:CD	2.14	0.51
23:DA:1973:G:H2'	23:DA:1974:C:H6	1.75	0.51
23:BA:2758:A:C4	29:BG:67:LEU:HD21	2.45	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DA:2188:C:H2'	23:DA:2189:U:O4'	2.10	0.51
25:BC:166:GLN:CA	25:BC:166:GLN:HE21	2.22	0.51
23:BA:566:U:H2'	23:BA:567:A:O4'	2.10	0.51
23:DA:1444:G:H2'	23:DA:1445:C:C5	2.46	0.51
23:DA:144(B):A:H5''	23:DA:1445:C:H5	1.75	0.51
23:DA:2032:G:O2'	26:DD:145:LYS:HE2	2.09	0.51
11:AK:73:MET:HG2	11:AK:103:LEU:HD11	1.91	0.51
23:DA:1028:A:N6	23:DA:1125:G:H2'	2.25	0.51
23:BA:830:G:H4'	23:BA:831:G:OP2	2.11	0.51
23:BA:780:G:N2	23:BA:783:A:H62	1.95	0.51
43:BU:37:VAL:HG21	43:BU:72:VAL:HG21	1.93	0.51
23:BA:848:G:C4	23:BA:933:A:C8	2.95	0.51
23:BA:2744:G:H21	29:BG:143:GLN:NE2	2.02	0.51
29:BG:92:ILE:CD1	29:BG:92:ILE:H	2.22	0.51
23:DA:2378:A:H2'	37:DO:21:THR:HG21	1.92	0.51
36:DN:21:TYR:HE2	36:DN:43:GLU:HB3	1.75	0.51
28:DF:74:LYS:HA	28:DF:74:LYS:HE3	1.92	0.51
24:DB:71:C:C2	24:DB:72:G:C8	2.98	0.51
27:BE:155:LEU:CD2	27:BE:186:ILE:HD13	2.39	0.51
37:BO:26:LEU:HD13	37:BO:87:PHE:HD1	1.76	0.51
30:DH:142:VAL:HG12	30:DH:143:SER:N	2.26	0.51
4:CD:189:PRO:CB	4:CD:194:LEU:HD21	2.38	0.51
23:BA:389:G:C6	34:BL:71:VAL:HG23	2.45	0.51
47:BY:46:GLN:HA	47:BY:46:GLN:OE1	2.09	0.51
9:AI:16:ARG:O	9:AI:63:ILE:HG23	2.10	0.51
9:CI:16:ARG:O	9:CI:63:ILE:HG23	2.10	0.51
1:AA:1520:G:H2'	1:AA:1521:G:C8	2.44	0.51
38:DP:6:LEU:O	38:DP:10:VAL:HG23	2.10	0.51
1:CA:983:A:H2	1:CA:984:C:C5	2.28	0.51
1:AA:1389:C:H2'	1:AA:1390:U:O4'	2.09	0.51
17:CQ:59:ILE:CG2	17:CQ:71:PHE:HB3	2.41	0.51
8:AH:11:THR:HA	8:AH:14:ARG:NH1	2.26	0.51
8:CH:80:ILE:N	8:CH:80:ILE:HD12	2.24	0.51
1:AA:1423:G:H2'	1:AA:1424:C:C6	2.45	0.51
2:CB:135:GLN:O	2:CB:139:LYS:HG2	2.10	0.51
23:DA:2252:G:H2'	23:DA:2253:G:H8	1.76	0.51
23:BA:943:U:OP2	34:BL:38:GLN:CD	2.48	0.51
23:BA:588:U:H1'	27:BE:90:PHE:CD1	2.45	0.51
23:BA:588:U:C2	27:BE:90:PHE:CE1	2.99	0.51
2:CB:69:LEU:HB3	2:CB:162:ILE:HG22	1.93	0.51
42:BT:30:VAL:HG11	42:BT:39:ILE:CD1	2.38	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
47:DY:12:GLU:C	47:DY:14:ARG:H	2.13	0.51
8:CH:66:GLY:HA3	8:CH:77:GLU:HB3	1.92	0.51
1:AA:1308:U:H5''	13:AM:98:VAL:HG23	1.92	0.51
1:CA:80:G:P	1:CA:80:G:H8	2.33	0.51
8:AH:103:VAL:HG21	8:AH:109:ILE:C	2.30	0.51
2:CB:27:LYS:HD3	2:CB:27:LYS:H	1.76	0.51
8:CH:103:VAL:HG21	8:CH:109:ILE:C	2.30	0.51
42:DT:15:GLU:CD	42:DT:15:GLU:N	2.62	0.51
28:DF:85:GLY:C	28:DF:86:MET:HG3	2.31	0.51
23:BA:661:C:H2'	23:BA:662:G:C8	2.46	0.51
4:CD:21:LEU:HD12	4:CD:22:LYS:H	1.75	0.51
30:DH:116:LEU:HD22	30:DH:128:LEU:HD21	1.91	0.51
30:DH:86:THR:O	30:DH:122:GLU:HG3	2.11	0.51
9:CI:113:LYS:HG2	9:CI:119:ALA:HA	1.92	0.51
27:BE:34:TRP:HB2	34:BL:10:PRO:O	2.10	0.51
1:AA:626:U:H2'	1:AA:627:G:C8	2.45	0.51
1:AA:1267:C:O2	1:AA:1327:C:H4'	2.10	0.51
23:BA:2105:C:H2'	23:BA:2106:G:C8	2.45	0.51
1:AA:59:A:H3'	1:AA:331:G:H22	1.75	0.51
37:BO:15:ARG:O	37:BO:19:LYS:HG3	2.10	0.51
15:AO:15:PHE:O	15:AO:27:VAL:HG22	2.10	0.51
22:AV:6166:U:C4	22:AV:6167:G:N7	2.78	0.51
15:CO:15:PHE:O	15:CO:27:VAL:HG22	2.10	0.51
7:CG:38:LEU:HD12	7:CG:41:ARG:HH12	1.76	0.51
1:CA:176:C:H5''	20:CT:29:LYS:NZ	2.25	0.51
3:AC:184:TYR:CE2	3:AC:186:PHE:HB2	2.45	0.51
1:CA:643:C:H5'	8:CH:31:PHE:CD1	2.45	0.51
23:BA:1444:G:H2'	23:BA:1445:C:C5	2.46	0.51
39:DQ:8:VAL:HG11	39:DQ:12:ARG:CZ	2.39	0.51
1:CA:164:U:H2'	1:CA:165:C:C6	2.45	0.51
23:DA:1010:A:H1'	23:DA:1153:C:H1'	1.93	0.51
39:BQ:107:ALA:O	39:BQ:111:GLU:HG2	2.11	0.51
23:BA:69:C:O2'	23:BA:70:G:H5'	2.11	0.51
23:BA:144:C:H2'	23:BA:145:G:C8	2.45	0.51
8:CH:11:THR:HG22	8:CH:15:ASN:ND2	2.26	0.51
20:AT:14:LYS:O	20:AT:18:GLN:HG3	2.10	0.51
1:CA:37:U:OP1	12:CL:122:LYS:HG3	2.10	0.51
41:BS:19:LEU:HB3	50:B2:25:LEU:CD1	2.40	0.51
39:BQ:106:PHE:O	39:BQ:110:VAL:HG23	2.10	0.51
34:DL:85:LEU:HD23	34:DL:85:LEU:H	1.75	0.51
34:BL:83:VAL:O	34:BL:114:ILE:HA	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BA:954:G:C5	23:BA:955:C:C5	2.99	0.51
23:BA:960:A:H61	35:BM:82:ARG:HH21	1.58	0.51
1:CA:78:G:H2'	1:CA:79:G:C8	2.44	0.51
13:CM:60:VAL:HG13	13:CM:64:TRP:NE1	2.22	0.51
3:AC:17:ASP:CB	3:AC:21:ARG:HH22	2.23	0.51
16:CP:4:ILE:HD13	16:CP:66:PRO:HG3	1.92	0.51
23:BA:603:A:H61	23:BA:655:A:C1'	2.24	0.51
25:BC:118:VAL:HG22	25:BC:119:ALA:N	2.24	0.51
36:BN:10:LEU:HB2	36:BN:17:ARG:CZ	2.41	0.51
23:BA:1188:U:C2'	23:BA:1189:A:H5'	2.40	0.51
23:DA:310:A:OP1	43:DU:18:GLY:HA2	2.11	0.51
28:BF:32:PRO:HA	28:BF:162:THR:OG1	2.10	0.51
18:AR:74:ARG:HA	18:AR:79:LEU:O	2.11	0.51
23:DA:1541:U:H3'	23:DA:1542:G:H3'	1.93	0.51
27:DE:182:ASN:O	27:DE:186:ILE:HG12	2.11	0.51
46:DX:27:GLU:HB2	46:DX:32:LYS:O	2.11	0.51
30:DH:101:LEU:HG	30:DH:107:ILE:HG23	1.92	0.51
4:AD:108:LEU:HB3	4:AD:110:PHE:CD2	2.45	0.51
4:CD:108:LEU:HD23	4:CD:110:PHE:CE2	2.44	0.51
23:BA:814:C:H5	34:BL:27:HIS:NE2	2.08	0.51
28:DF:134:GLY:C	28:DF:135:LEU:HD12	2.30	0.51
23:DA:320:A:H2'	27:DE:136:THR:HG21	1.91	0.51
35:BM:58:PHE:CD1	35:BM:61:GLY:HA3	2.45	0.51
22:AV:6168:G:H2'	22:AV:6169:U:C6	2.46	0.51
23:BA:719:C:H2'	23:BA:720:C:H6	1.75	0.51
5:CE:72:GLN:O	5:CE:75:THR:HG22	2.09	0.51
23:DA:915:C:H2'	23:DA:916:G:C8	2.46	0.51
23:BA:1824:G:OP1	25:BC:52:ARG:HD3	2.11	0.51
8:AH:31:PHE:O	8:AH:35:ILE:HG12	2.10	0.51
15:CO:26:GLU:OE2	15:CO:77:ARG:HD2	2.11	0.51
23:DA:1206:G:C6	23:DA:1207:C:C4	2.98	0.51
23:DA:498:G:N3	43:DU:47:LYS:HE3	2.25	0.51
4:AD:142:PRO:HA	4:AD:185:PHE:HD2	1.74	0.51
23:BA:1344:G:H4'	23:BA:1384:A:C5	2.46	0.51
30:BH:31:LEU:HB3	30:BH:32:PRO:HD3	1.92	0.51
1:AA:721:G:H4'	1:AA:722:A:O4'	2.11	0.51
36:BN:72:ASP:O	36:BN:76:VAL:HG12	2.11	0.51
5:AE:76:ILE:HD11	5:AE:142:LEU:HD11	1.91	0.51
23:BA:1022:G:H8	32:BJ:92:GLN:HE22	1.57	0.51
35:DM:75:THR:CA	35:DM:88:GLY:HA2	2.39	0.51
46:DX:11:ARG:HB2	46:DX:13:ILE:HG22	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
29:BG:92:ILE:N	29:BG:92:ILE:HD12	2.22	0.51
25:BC:71:ASP:CB	25:BC:103:ARG:HH22	2.20	0.51
1:AA:715:A:H2'	1:AA:716:A:C8	2.46	0.51
23:DA:1335:U:H2'	23:DA:1336:A:H8	1.76	0.51
23:DA:389:G:C6	34:DL:71:VAL:HG23	2.45	0.51
15:CO:45:VAL:HG23	15:CO:46:HIS:ND1	2.26	0.51
23:DA:1187:G:H8	23:DA:1187:G:O5'	1.94	0.51
1:CA:434:U:H2'	1:CA:435:C:H6	1.74	0.51
22:AV:6168:G:H2'	22:AV:6169:U:H6	1.76	0.51
22:CV:6166:U:C4	22:CV:6167:G:N7	2.78	0.51
23:DA:2537:U:H2'	23:DA:2538:C:H6	1.75	0.51
34:DL:75:ILE:HD13	34:DL:77:ARG:NE	2.25	0.51
23:BA:363(D):G:H2'	23:BA:363(E):G:H8	1.76	0.51
6:CF:21:LEU:O	6:CF:25:ILE:HG12	2.10	0.51
4:AD:100:ARG:NH1	4:AD:137:SER:HA	2.24	0.51
15:AO:26:GLU:OE2	15:AO:77:ARG:HD2	2.10	0.51
37:BO:49:VAL:HG11	37:BO:73:LEU:HA	1.92	0.51
1:AA:321:A:H2'	1:AA:322:C:C6	2.46	0.51
3:AC:58:GLU:HB2	3:AC:65:ALA:HB3	1.92	0.51
23:BA:2378:A:H2'	37:BO:21:THR:HG21	1.92	0.51
25:DC:244:ARG:HG3	25:DC:245:PRO:N	2.25	0.51
24:BB:71:C:C4	24:BB:72:G:N7	2.79	0.51
28:BF:77:ILE:HG22	28:BF:80:PHE:N	2.24	0.51
23:DA:588:U:C2	27:DE:90:PHE:CE1	2.98	0.51
22:AV:6157:A:N6	22:AV:6172:U:H3	2.09	0.51
23:BA:2893:G:H3'	23:BA:2894:G:H5'	1.93	0.51
23:DA:2893:G:H3'	23:DA:2894:G:H5'	1.93	0.51
1:CA:1520:G:H2'	1:CA:1521:G:C8	2.46	0.51
1:AA:1190:G:OP2	3:AC:5:ILE:HG23	2.11	0.51
39:BQ:49:HIS:HA	39:BQ:52:ARG:HB2	1.91	0.51
23:BA:581:C:OP1	39:BQ:31:SER:HB2	2.11	0.51
23:BA:920:G:H2'	23:BA:921:G:H8	1.76	0.51
11:AK:99:GLN:HE22	11:AK:105:VAL:HG21	1.76	0.51
37:DO:38:GLN:HB3	37:DO:47:THR:CG2	2.40	0.51
1:AA:1203:C:H2'	1:AA:1204:A:H8	1.76	0.51
23:DA:23:G:H2'	23:DA:24:G:H8	1.75	0.51
1:CA:1290:G:H2'	1:CA:1290:G:N3	2.26	0.51
15:AO:45:VAL:HG23	15:AO:46:HIS:ND1	2.25	0.51
40:BR:13:ARG:C	40:BR:13:ARG:HD2	2.31	0.51
48:BZ:11:SER:OG	48:BZ:13:ILE:HG12	2.10	0.51
52:D4:24:THR:O	52:D4:28:ARG:HG3	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
39:BQ:90:VAL:HG13	39:BQ:91:ASP:N	2.26	0.51
39:BQ:95:LEU:HD11	40:BR:12:TYR:HA	1.92	0.51
47:DY:14:ARG:HA	47:DY:17:SER:HB2	1.93	0.51
53:D5:50:LEU:HD13	53:D5:57:ARG:CZ	2.40	0.51
13:AM:60:VAL:HG13	13:AM:64:TRP:NE1	2.23	0.51
43:BU:71:LYS:HB2	43:BU:71:LYS:HZ3	1.76	0.51
23:DA:1404:C:O2'	23:DA:1405:U:H5'	2.10	0.51
23:DA:917:A:H2'	23:DA:918:A:O4'	2.10	0.51
19:AS:22:LEU:HD13	19:AS:27:GLU:HB2	1.93	0.51
6:CF:69:GLU:H	6:CF:69:GLU:CD	2.15	0.51
44:BV:104:PHE:HB3	44:BV:141:VAL:HG11	1.93	0.51
23:BA:388:G:OP1	46:BX:33:LYS:HB3	2.11	0.51
4:CD:122:ARG:C	4:CD:122:ARG:HD3	2.32	0.51
20:CT:26:ASN:HD22	20:CT:27:LYS:N	2.07	0.51
23:BA:2777:G:C5'	23:BA:2778:A:H5'	2.41	0.51
42:DT:89:ILE:O	42:DT:93:GLU:HG2	2.11	0.51
34:BL:135:LEU:O	34:BL:135:LEU:HD13	2.11	0.51
23:BA:379:G:N2	46:BX:20:ARG:HH12	2.08	0.51
23:BA:2698:U:H2'	23:BA:2699:C:C6	2.46	0.51
1:CA:161:A:H2'	1:CA:162:A:C8	2.46	0.51
1:CA:501:C:H2'	1:CA:502:G:H8	1.74	0.51
1:AA:189:U:O4	17:AQ:62:SER:HB3	2.11	0.51
1:CA:10:A:H2'	1:CA:11:G:H8	1.76	0.51
23:DA:1817:G:OP1	25:DC:88:ARG:NH2	2.43	0.51
23:DA:49:A:H5''	23:DA:51:G:O4'	2.10	0.51
34:DL:138:LEU:HD11	34:DL:144:GLU:HB3	1.93	0.51
23:DA:719:C:H2'	23:DA:720:C:H6	1.75	0.51
35:DM:74:TYR:CD2	35:DM:91:GLU:HB2	2.46	0.51
2:AB:135:GLN:O	2:AB:139:LYS:HG2	2.11	0.51
1:AA:824:C:H3'	1:AA:824:C:C6	2.46	0.51
17:CQ:52:LYS:H	17:CQ:52:LYS:HD2	1.76	0.51
1:AA:617:G:H4'	16:AP:44:THR:HB	1.93	0.51
44:BV:118:GLN:HB2	44:BV:173:ALA:O	2.11	0.51
16:CP:23:ASP:O	16:CP:26:ARG:HB2	2.10	0.51
36:BN:104:ARG:NH1	36:BN:109:ALA:HB3	2.25	0.51
9:CI:19:LEU:HD23	9:CI:20:ARG:N	2.26	0.51
2:CB:163:PHE:HD1	2:CB:185:ILE:HG13	1.75	0.51
8:AH:66:GLY:HA3	8:AH:77:GLU:HB3	1.92	0.51
1:AA:89:U:H2'	1:AA:90:C:C6	2.46	0.51
53:D5:53:PRO:HB2	53:D5:57:ARG:NH2	2.26	0.51
12:CL:31:PHE:HB3	12:CL:83:LEU:HD22	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:DN:17:ARG:O	36:DN:20:LEU:HB3	2.11	0.51
19:CS:22:LEU:HD13	19:CS:27:GLU:HB2	1.93	0.51
34:BL:16:ARG:HE	34:BL:16:ARG:C	2.14	0.51
25:DC:118:VAL:HG22	25:DC:119:ALA:N	2.25	0.51
37:DO:26:LEU:HG	37:DO:39:ILE:CD1	2.40	0.51
1:AA:1225:A:H5''	1:AA:1226:C:OP2	2.11	0.51
30:DH:4:ILE:HD11	30:DH:16:GLY:HA2	1.92	0.51
1:AA:1152:A:H2'	1:AA:1153:C:C6	2.46	0.51
23:BA:597:U:O2'	34:BL:15:ARG:HG2	2.10	0.51
2:CB:178:ARG:HD2	8:CH:71:GLY:O	2.11	0.51
38:BP:100:TYR:HD2	38:BP:103:ARG:NH2	2.09	0.51
3:CC:57:ILE:HD13	3:CC:66:VAL:HA	1.93	0.51
24:DB:30:C:H2'	24:DB:31:C:H5'	1.93	0.51
23:BA:2305:A:H3'	23:BA:2306:C:H5''	1.92	0.51
7:AG:38:LEU:HD12	7:AG:41:ARG:HH12	1.76	0.51
23:DA:2698:U:H2'	23:DA:2699:C:C6	2.46	0.51
21:AU:6:ARG:NE	21:AU:15:ARG:HH12	2.09	0.51
35:DM:68:ILE:HG23	35:DM:103:MET:HA	1.92	0.51
15:CO:53:HIS:HE1	23:DA:715:G:O6	1.94	0.51
23:BA:2795:G:H3'	23:BA:2797:U:C5'	2.41	0.51
17:CQ:55:ASP:HB3	17:CQ:76:LEU:HD13	1.91	0.51
8:CH:110:ALA:HB3	8:CH:121:ASP:HB3	1.93	0.51
51:B3:30:THR:HG22	51:B3:31:PRO:HD2	1.92	0.51
23:DA:2795:G:H3'	23:DA:2797:U:C5'	2.41	0.51
35:BM:74:TYR:CD2	35:BM:91:GLU:HB2	2.46	0.51
34:DL:147:LEU:HD13	34:DL:148:LEU:O	2.11	0.51
43:DU:95:LYS:HG2	43:DU:100:ALA:HA	1.92	0.51
35:DM:75:THR:HG21	35:DM:85:LYS:HZ1	1.74	0.51
11:AK:29:ILE:C	11:AK:29:ILE:HD12	2.31	0.51
30:DH:72:LEU:HD12	30:DH:140:LEU:HD13	1.92	0.51
30:BH:82:ARG:HG2	30:BH:89:TYR:CD1	2.46	0.51
23:BA:1335:U:H2'	23:BA:1336:A:H8	1.76	0.51
13:AM:3:ARG:HH21	13:AM:7:VAL:HG13	1.75	0.51
23:BA:2484:G:H5''	35:BM:45:GLN:HB2	1.92	0.51
12:CL:65:VAL:HG11	12:CL:97:TYR:CE1	2.45	0.51
33:DK:2:ILE:CD1	33:DK:82:ASN:HD22	2.23	0.51
1:CA:475:G:H2'	1:CA:476:G:C8	2.46	0.51
25:BC:43:ARG:HB2	25:BC:48:ARG:O	2.11	0.51
41:BS:46:PHE:O	41:BS:50:VAL:HG12	2.11	0.51
2:CB:96:ARG:N	2:CB:96:ARG:HD2	2.25	0.51
35:DM:48:GLU:O	35:DM:52:VAL:HG12	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
17:CQ:59:ILE:HD12	17:CQ:59:ILE:N	2.26	0.51
23:BA:144(B):A:H5''	23:BA:1445:C:H5	1.75	0.51
39:DQ:8:VAL:HG11	39:DQ:12:ARG:NE	2.26	0.51
23:BA:2840:C:H4'	36:BN:53:HIS:CD2	2.46	0.51
11:CK:105:VAL:HG23	11:CK:105:VAL:O	2.11	0.51
11:CK:73:MET:HG2	11:CK:103:LEU:HD11	1.92	0.51
25:DC:37:LEU:HD12	25:DC:38:LYS:H	1.75	0.51
23:BA:2853:C:H2'	23:BA:2854:G:H8	1.76	0.51
1:AA:836:G:C6	1:AA:851:G:C6	2.99	0.51
36:DN:72:ASP:O	36:DN:76:VAL:HG12	2.10	0.51
23:DA:1809:A:H2'	23:DA:1810:A:C8	2.46	0.51
17:CQ:74:LEU:HD12	17:CQ:75:ARG:HG2	1.93	0.51
27:BE:184:TYR:CE2	27:BE:188:ARG:HD2	2.46	0.51
23:DA:566:U:H2'	23:DA:567:A:O4'	2.11	0.51
40:DR:13:ARG:HD2	40:DR:13:ARG:C	2.31	0.51
3:AC:79:ARG:HD3	3:AC:79:ARG:N	2.26	0.51
1:CA:963:G:H2'	1:CA:964:A:C8	2.46	0.51
24:DB:63:G:H2'	24:DB:64:C:C6	2.46	0.51
23:DA:1680:U:O2	23:DA:1763:G:H3'	2.11	0.51
25:DC:238:GLY:O	25:DC:240:ALA:N	2.44	0.50
34:BL:85:LEU:HD23	34:BL:85:LEU:H	1.76	0.50
23:DA:955:C:OP2	35:DM:14:ARG:HD2	2.11	0.50
35:DM:8:LYS:O	35:DM:9:TYR:HB3	2.11	0.50
42:DT:30:VAL:HG11	42:DT:39:ILE:CD1	2.38	0.50
46:DX:62:VAL:HG22	46:DX:63:ALA:N	2.26	0.50
1:AA:674:G:H2'	1:AA:675:A:C8	2.46	0.50
25:DC:131:LEU:HD11	25:DC:136:ILE:HG12	1.92	0.50
1:CA:38:G:C2	1:CA:397:A:C2	2.99	0.50
23:BA:744:G:OP1	26:BD:132:HIS:HB2	2.12	0.50
1:CA:1014:A:H2	1:CA:1219:U:H1'	1.75	0.50
23:BA:2038:G:H2'	23:BA:2039:C:H6	1.76	0.50
2:AB:27:LYS:H	2:AB:27:LYS:HD3	1.75	0.50
23:BA:94:G:H21	47:BY:47:ASN:ND2	2.09	0.50
18:AR:36:ASN:HD22	18:AR:39:VAL:HG21	1.76	0.50
44:BV:5:LEU:HB3	44:BV:59:LEU:HD23	1.92	0.50
23:DA:628:G:H2'	23:DA:629:G:H8	1.75	0.50
1:AA:983:A:H3'	1:AA:983:A:N3	2.26	0.50
23:DA:2419:U:O4	53:D5:30:ARG:CZ	2.60	0.50
23:DA:1537:C:H2'	23:DA:1538:G:O4'	2.11	0.50
23:DA:1178:C:H2'	23:DA:1179:C:H6	1.75	0.50
1:CA:1260:C:H4'	1:CA:1284:C:H5'	1.92	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:772:U:H2'	1:AA:773:G:O4'	2.10	0.50
1:CA:127:G:HO2'	17:CQ:2:PRO:N	2.09	0.50
23:DA:1929:G:H5''	23:DA:1929:G:N3	2.26	0.50
23:BA:2366:A:H2'	23:BA:2367:G:O4'	2.11	0.50
34:BL:33:ARG:HG3	34:BL:36:LYS:CD	2.22	0.50
23:DA:960:A:H61	35:DM:82:ARG:HH21	1.60	0.50
23:DA:675:A:H4'	27:DE:67:GLN:HE21	1.76	0.50
1:AA:979:C:H42	14:AN:18:VAL:HG12	1.76	0.50
23:BA:2822:G:H8	23:BA:2822:G:O5'	1.95	0.50
23:DA:390:A:C6	34:DL:71:VAL:HG21	2.46	0.50
23:DA:646:A:N3	23:DA:646:A:H5'	2.27	0.50
43:DU:20:TYR:CE1	43:DU:42:VAL:HA	2.46	0.50
20:CT:67:ALA:HA	20:CT:72:LEU:O	2.12	0.50
46:BX:90:ILE:O	46:BX:94:LEU:HB2	2.10	0.50
23:DA:2777:G:C5'	23:DA:2778:A:H5'	2.42	0.50
32:DJ:127:LYS:HA	32:DJ:130:LEU:HD12	1.93	0.50
3:AC:57:ILE:HD13	3:AC:66:VAL:HA	1.94	0.50
1:CA:464:G:O5'	1:CA:464:G:H8	1.94	0.50
23:DA:814:C:H5	34:DL:27:HIS:NE2	2.09	0.50
46:BX:45:ASN:ND2	46:BX:47:GLN:HE21	2.10	0.50
1:AA:665:A:H2'	1:AA:732:C:O2	2.11	0.50
23:DA:150:C:H2'	23:DA:151:C:C6	2.46	0.50
1:AA:1306:A:H2'	1:AA:1307:U:C6	2.46	0.50
23:BA:1111:A:N3	23:BA:1112:G:H1'	2.26	0.50
1:CA:1389:C:H2'	1:CA:1390:U:O4'	2.10	0.50
1:CA:447:G:H2'	1:CA:485:G:N2	2.26	0.50
28:DF:133:LEU:HD21	28:DF:157:ILE:HG13	1.93	0.50
53:B5:11:LYS:HD3	53:B5:11:LYS:C	2.31	0.50
8:CH:81:HIS:HB2	8:CH:138:TRP:OXT	2.11	0.50
1:CA:1201:A:H4'	1:CA:1202:G:O5'	2.10	0.50
23:BA:2340:G:H2'	23:BA:2341:G:H8	1.76	0.50
37:BO:36:TYR:CD1	37:BO:36:TYR:N	2.79	0.50
27:DE:179:GLU:CD	27:DE:179:GLU:H	2.13	0.50
1:CA:1412:C:H2'	1:CA:1413:A:C8	2.47	0.50
23:BA:593:G:O2'	53:B5:62:LEU:HD13	2.11	0.50
39:DQ:112:ARG:HH21	40:DR:46:VAL:HG21	1.76	0.50
39:DQ:88:ILE:HB	39:DQ:90:VAL:CG1	2.35	0.50
39:BQ:112:ARG:HH21	40:BR:46:VAL:HG21	1.76	0.50
37:BO:30:ARG:C	37:BO:30:ARG:HD2	2.31	0.50
46:BX:62:VAL:HG22	46:BX:63:ALA:N	2.26	0.50
10:CJ:75:ILE:HG13	10:CJ:76:ASN:N	2.22	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DA:1813:G:H1'	25:DC:50:THR:CG2	2.37	0.50
25:BC:33:LEU:HD23	25:BC:33:LEU:N	2.24	0.50
12:CL:81:VAL:O	12:CL:82:VAL:HB	2.12	0.50
42:BT:63:LYS:HZ2	42:BT:72:LYS:HB3	1.76	0.50
8:AH:6:ILE:O	8:AH:10:LEU:HG	2.12	0.50
37:BO:26:LEU:HG	37:BO:39:ILE:CD1	2.41	0.50
37:DO:26:LEU:HD13	37:DO:87:PHE:HD1	1.76	0.50
23:BA:1541:U:H3'	23:BA:1542:G:H3'	1.93	0.50
12:AL:44:PRO:HD2	12:AL:49:SER:HA	1.93	0.50
30:DH:107:ILE:HG13	30:DH:109:ILE:HG23	1.92	0.50
23:BA:1404:C:O2'	23:BA:1405:U:H5'	2.11	0.50
23:BA:1504:C:O2'	23:BA:1505:C:H6	1.95	0.50
1:AA:1148:U:C2	9:AI:16:ARG:NH2	2.79	0.50
23:BA:2271:G:H2'	23:BA:2272:U:C6	2.46	0.50
42:BT:89:ILE:O	42:BT:93:GLU:HG2	2.11	0.50
38:DP:57:PHE:CG	38:DP:58:ASN:N	2.80	0.50
44:DV:39:VAL:HG21	44:DV:44:PHE:CD2	2.47	0.50
8:CH:11:THR:HA	8:CH:14:ARG:NH1	2.26	0.50
11:CK:99:GLN:HE22	11:CK:105:VAL:HG21	1.76	0.50
23:BA:1178:C:H2'	23:BA:1179:C:H6	1.76	0.50
23:DA:2641:G:H5''	32:DJ:99:SER:HB3	1.94	0.50
9:AI:4:TYR:CE2	9:AI:88:TYR:HB2	2.47	0.50
23:BA:2208:U:O2'	23:BA:2209:C:H5'	2.11	0.50
23:DA:1165:U:H2'	23:DA:1166:C:C6	2.47	0.50
44:DV:27:VAL:HG22	44:DV:36:LYS:HA	1.93	0.50
26:BD:172:VAL:HG13	26:BD:182:LEU:HD11	1.92	0.50
32:DJ:49:LEU:O	32:DJ:53:ILE:HG13	2.11	0.50
52:B4:24:THR:O	52:B4:28:ARG:HG3	2.11	0.50
10:AJ:29:ARG:HG2	10:AJ:29:ARG:O	2.12	0.50
23:BA:1929:G:H5''	23:BA:1929:G:N3	2.26	0.50
26:BD:14:ILE:C	26:BD:14:ILE:HD12	2.31	0.50
2:AB:96:ARG:N	2:AB:96:ARG:HD2	2.25	0.50
23:DA:2335:A:C8	23:DA:2337:G:C5	3.00	0.50
23:DA:379:G:N2	46:DX:20:ARG:HH12	2.09	0.50
34:BL:147:LEU:HD13	34:BL:148:LEU:O	2.12	0.50
23:DA:780:G:N2	23:DA:783:A:H62	1.96	0.50
2:AB:69:LEU:HB3	2:AB:162:ILE:HG22	1.92	0.50
37:DO:30:ARG:HD2	37:DO:30:ARG:C	2.32	0.50
44:BV:125:LEU:HD23	44:BV:126:VAL:N	2.26	0.50
25:BC:244:ARG:HH11	25:BC:244:ARG:HB2	1.77	0.50
12:AL:81:VAL:O	12:AL:82:VAL:HB	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:AB:24:TRP:HE3	2:AB:25:ASN:O	1.95	0.50
32:BJ:57:LEU:HD11	32:BJ:142:ARG:HB2	1.92	0.50
28:DF:74:LYS:HE2	28:DF:84:LYS:HE3	1.94	0.50
11:CK:20:TYR:O	11:CK:30:VAL:HA	2.11	0.50
11:CK:21:ILE:HD12	11:CK:21:ILE:N	2.26	0.50
37:DO:25:ARG:HD2	37:DO:88:ASP:OD1	2.12	0.50
25:DC:43:ARG:HB2	25:DC:48:ARG:O	2.10	0.50
1:AA:1356:G:H2'	1:AA:1357:A:H8	1.74	0.50
4:CD:105:VAL:HG13	4:CD:110:PHE:HB2	1.93	0.50
4:CD:108:LEU:HB3	4:CD:110:PHE:CD2	2.46	0.50
32:BJ:127:LYS:HA	32:BJ:130:LEU:HD12	1.92	0.50
1:AA:841:U:HO2'	1:AA:842:C:H6	1.59	0.50
27:BE:12:LEU:HB2	27:BE:124:LEU:HD11	1.94	0.50
23:DA:628:G:H2'	23:DA:629:G:C8	2.47	0.50
17:CQ:57:VAL:HG12	17:CQ:76:LEU:HA	1.94	0.50
1:AA:983:A:H2	1:AA:984:C:C5	2.29	0.50
23:DA:2462:U:H1'	23:DA:2491:U:O4	2.10	0.50
23:BA:1010:A:H1'	23:BA:1153:C:H1'	1.92	0.50
23:DA:576:U:H2'	23:DA:577:G:C8	2.46	0.50
2:CB:75:LYS:C	2:CB:75:LYS:HD3	2.32	0.50
23:DA:1056:G:H8	23:DA:1056:G:O5'	1.94	0.50
23:BA:2462:U:H1'	23:BA:2491:U:O4	2.12	0.50
7:CG:113:GLU:HB3	7:CG:118:VAL:HG23	1.93	0.50
39:DQ:90:VAL:HG13	39:DQ:91:ASP:N	2.26	0.50
40:BR:49:THR:HB	40:BR:50:PRO:CD	2.42	0.50
43:DU:14:LEU:HD23	43:DU:15:VAL:N	2.26	0.50
42:DT:11:PRO:HG3	47:DY:37:PHE:CE2	2.46	0.50
1:CA:90:C:H2'	1:CA:91:C:C6	2.47	0.50
44:DV:125:LEU:HD23	44:DV:126:VAL:N	2.27	0.50
1:AA:38:G:C2	1:AA:397:A:C2	3.00	0.50
1:CA:321:A:H2'	1:CA:322:C:C6	2.47	0.50
37:BO:14:VAL:O	37:BO:18:ILE:HG12	2.12	0.50
28:BF:74:LYS:HE2	28:BF:84:LYS:HE3	1.94	0.50
32:DJ:57:LEU:HD11	32:DJ:142:ARG:HB2	1.93	0.50
23:BA:2517:C:C5	23:BA:2542:A:C2	2.99	0.50
19:CS:32:LYS:HA	19:CS:50:ALA:HB3	1.94	0.50
5:CE:36:ASP:O	5:CE:37:ARG:HB2	2.12	0.50
28:BF:25:TYR:CD1	28:BF:30:GLU:HB3	2.47	0.50
29:DG:46:GLU:HG3	29:DG:51:ARG:NE	2.26	0.50
18:CR:74:ARG:HA	18:CR:79:LEU:O	2.11	0.50
1:AA:1072:G:H2'	1:AA:1073:U:C6	2.46	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BA:270(H):C:H2'	23:BA:270(I):C:H6	1.76	0.50
13:CM:49:THR:O	13:CM:53:VAL:HG23	2.12	0.50
30:DH:31:LEU:HB3	30:DH:32:PRO:HD3	1.94	0.50
1:CA:629:G:H2'	1:CA:630:G:C8	2.47	0.50
19:AS:53:ASN:HD21	19:AS:56:GLN:H	1.59	0.50
23:DA:2758:A:C4	29:DG:67:LEU:HD21	2.47	0.50
23:DA:363(D):G:H2'	23:DA:363(E):G:H8	1.77	0.50
23:BA:1056:G:O5'	23:BA:1056:G:H8	1.94	0.50
12:AL:19:LYS:H	12:AL:19:LYS:HD3	1.77	0.50
37:DO:36:TYR:N	37:DO:36:TYR:CD1	2.79	0.50
29:BG:30:LYS:HB2	29:BG:79:VAL:HA	1.94	0.50
1:AA:444:C:H2'	1:AA:445:G:C8	2.47	0.50
23:DA:1824:G:OP1	25:DC:52:ARG:HD3	2.11	0.50
4:AD:88:VAL:O	4:AD:92:VAL:HG23	2.12	0.50
27:DE:29:ASN:H	27:DE:112:MET:CE	2.25	0.50
37:DO:69:VAL:O	37:DO:72:ALA:HB3	2.11	0.50
25:BC:131:LEU:HD11	25:BC:136:ILE:HG12	1.94	0.50
29:DG:92:ILE:N	29:DG:92:ILE:HD12	2.21	0.50
25:DC:132:PRO:HG3	25:DC:190:TYR:CE1	2.46	0.50
36:DN:10:LEU:CB	36:DN:17:ARG:NE	2.73	0.50
2:CB:47:THR:O	2:CB:51:LEU:HG	2.12	0.50
23:DA:557:U:H2'	23:DA:558:G:H8	1.76	0.50
23:DA:518:G:H4'	41:DS:18:ARG:NH1	2.26	0.50
30:DH:6:LEU:HD23	30:DH:6:LEU:N	2.25	0.50
26:BD:120:TRP:CD1	26:BD:155:LYS:HB3	2.47	0.50
1:CA:370:C:H2'	1:CA:371:G:C8	2.47	0.50
23:DA:774:A:H2	23:DA:787:U:O2'	1.95	0.50
40:DR:81:TYR:O	40:DR:82:ARG:HG3	2.12	0.50
23:BA:775:G:C5	23:BA:794:G:C8	2.99	0.50
23:DA:1567:A:H2'	25:DC:84:TYR:HE2	1.76	0.50
23:BA:1980:G:H3'	23:BA:1981:A:C5'	2.42	0.50
1:AA:434:U:H2'	1:AA:435:C:H6	1.76	0.50
23:DA:2846:G:H2'	23:DA:2847:U:O4'	2.12	0.50
23:BA:116:C:H2'	23:BA:117:G:C8	2.47	0.50
23:BA:2320:A:H2'	23:BA:2320:A:N3	2.25	0.50
13:CM:17:VAL:HG12	13:CM:21:TYR:HE1	1.75	0.50
1:CA:824:C:H3'	1:CA:824:C:C6	2.47	0.50
2:CB:74:LYS:O	2:CB:78:GLN:HG3	2.12	0.50
6:CF:36:ARG:HH21	6:CF:38:GLU:HG2	1.75	0.50
39:DQ:95:LEU:HD11	40:DR:12:TYR:HA	1.93	0.50
43:BU:8:LYS:NZ	43:BU:8:LYS:N	2.57	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:922:G:N3	1:CA:1398:A:H2	2.10	0.50
10:AJ:33:GLN:O	10:AJ:75:ILE:HG12	2.12	0.50
42:DT:63:LYS:HZ2	42:DT:72:LYS:HB3	1.77	0.50
1:AA:922:G:N3	1:AA:1398:A:H2	2.09	0.50
3:CC:30:ARG:CD	14:CN:38:GLY:HA3	2.40	0.50
23:DA:1405:U:H2'	23:DA:1406:U:H6	1.75	0.50
30:BH:77:LEU:HG	30:BH:101:LEU:HD13	1.94	0.50
34:BL:17:LYS:O	34:BL:19:VAL:HG22	2.12	0.50
28:DF:25:TYR:CD1	28:DF:30:GLU:HB3	2.47	0.50
22:AV:6171:U:H2'	22:AV:6172:U:C6	2.47	0.50
1:CA:273:A:H1'	17:CQ:16:GLN:OE1	2.11	0.50
23:BA:481:G:H1'	23:BA:506:G:N2	2.27	0.50
2:CB:7:VAL:O	2:CB:11:LEU:HG	2.12	0.50
2:AB:7:VAL:O	2:AB:11:LEU:HG	2.12	0.50
30:BH:6:LEU:HD23	30:BH:6:LEU:N	2.26	0.50
25:DC:150:LYS:HE3	25:DC:150:LYS:HA	1.94	0.50
2:AB:178:ARG:HD2	8:AH:71:GLY:O	2.12	0.50
23:DA:1602:U:H3'	23:DA:1603:A:H5''	1.94	0.50
3:CC:91:LEU:HB3	3:CC:99:VAL:HG11	1.94	0.50
38:BP:57:PHE:CG	38:BP:58:ASN:N	2.80	0.50
23:DA:1504:C:O2'	23:DA:1505:C:H6	1.95	0.50
23:DA:2244:U:O2'	23:DA:2245:U:H5'	2.11	0.50
41:BS:4:LYS:HG2	41:BS:106:ILE:CG2	2.42	0.50
23:BA:1495:A:N3	23:BA:1495:A:H2'	2.26	0.50
5:CE:12:LEU:HD22	5:CE:12:LEU:C	2.32	0.50
20:CT:10:LEU:O	20:CT:13:LEU:HD13	2.11	0.50
27:DE:28:ILE:O	27:DE:30:PRO:HD3	2.11	0.50
29:DG:23:ARG:N	29:DG:23:ARG:HD3	2.27	0.50
17:AQ:74:LEU:HD12	17:AQ:75:ARG:HG2	1.93	0.50
1:AA:447:G:H2'	1:AA:485:G:N2	2.27	0.50
1:AA:1123:A:H4'	10:AJ:36:GLY:HA3	1.94	0.50
37:BO:38:GLN:HB3	37:BO:47:THR:CG2	2.41	0.50
28:DF:16:ARG:O	28:DF:20:ILE:HG12	2.12	0.50
45:BW:37:LEU:HG	45:BW:60:PHE:HA	1.92	0.50
39:DQ:106:PHE:O	39:DQ:110:VAL:HG23	2.12	0.50
23:DA:2246:G:H2'	23:DA:2247:A:C8	2.47	0.50
34:DL:83:VAL:O	34:DL:114:ILE:HA	2.11	0.50
1:AA:90:C:H2'	1:AA:91:C:C6	2.47	0.50
23:BA:310:A:OP1	43:BU:18:GLY:HA2	2.11	0.50
8:CH:6:ILE:O	8:CH:10:LEU:HG	2.11	0.50
4:CD:121:VAL:O	4:CD:134:ASP:HA	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
20:CT:26:ASN:HD22	20:CT:27:LYS:H	1.60	0.50
23:DA:37:C:H2'	23:DA:38:A:C8	2.47	0.50
1:AA:464:G:O5'	1:AA:464:G:H8	1.94	0.50
12:AL:40:ARG:HD3	12:AL:41:THR:O	2.12	0.50
2:CB:17:PHE:CD1	2:CB:44:LEU:HD21	2.47	0.50
10:AJ:30:SER:HB2	10:AJ:80:LYS:HG2	1.93	0.50
1:CA:909:A:H2'	1:CA:910:C:O4'	2.11	0.50
23:DA:2572:A:H62	26:DD:145:LYS:HG3	1.77	0.50
23:BA:41:C:H2'	23:BA:43:G:O4'	2.11	0.50
24:BB:46:A:H2'	24:BB:47:C:C6	2.47	0.50
38:DP:107:ASP:O	38:DP:110:ILE:HG22	2.11	0.50
23:DA:1341:U:O4	42:DT:16:LYS:HE2	2.12	0.50
33:BK:61:VAL:HG13	33:BK:61:VAL:O	2.12	0.50
1:AA:350:G:O2'	1:AA:351:G:H5'	2.11	0.50
23:BA:2641:G:H5''	32:BJ:99:SER:HB3	1.93	0.50
1:AA:629:G:H2'	1:AA:630:G:C8	2.47	0.50
25:DC:81:ALA:HB3	25:DC:94:LEU:HB3	1.93	0.50
3:CC:195:VAL:CG1	3:CC:196:LEU:H	2.20	0.50
23:BA:773:U:H4'	25:BC:47:GLY:CA	2.39	0.50
3:CC:58:GLU:HB2	3:CC:65:ALA:HB3	1.92	0.50
23:BA:646:A:N3	23:BA:646:A:H5'	2.27	0.50
28:BF:173:LEU:HA	28:BF:176:LEU:HD12	1.94	0.50
23:DA:860:U:O2	23:DA:860:U:O4'	2.30	0.50
1:CA:1072:G:H2'	1:CA:1073:U:C6	2.47	0.50
1:AA:1080:A:H5''	1:AA:1081:G:OP2	2.12	0.50
1:CA:1070:U:H2'	1:CA:1071:C:C6	2.47	0.50
30:BH:4:ILE:HD11	30:BH:16:GLY:HA2	1.94	0.50
38:DP:100:TYR:HD2	38:DP:103:ARG:NH2	2.10	0.50
23:BA:37:C:H2'	23:BA:38:A:C8	2.46	0.50
23:DA:1603:A:H8	23:DA:1603:A:H5'	1.77	0.50
38:DP:54:ARG:HA	38:DP:59:THR:OG1	2.12	0.50
25:BC:150:LYS:HA	25:BC:150:LYS:HE3	1.94	0.50
28:DF:130:ASN:OD1	28:DF:160:VAL:HA	2.11	0.50
44:BV:39:VAL:HG21	44:BV:44:PHE:CD2	2.47	0.50
20:CT:85:MET:HB2	20:CT:104:LEU:HD21	1.93	0.50
2:AB:17:PHE:CD1	2:AB:44:LEU:HD21	2.47	0.50
23:BA:2846:G:H2'	23:BA:2847:U:O4'	2.11	0.50
19:CS:12:ASP:HB2	19:CS:15:LEU:HD23	1.93	0.50
23:DA:1344:G:H4'	23:DA:1384:A:C5	2.47	0.50
43:DU:29:GLU:HB3	43:DU:38:ILE:HB	1.94	0.50
5:AE:65:ASN:O	5:AE:66:MET:HB2	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:CD:162:LEU:HD13	4:CD:181:MET:HG2	1.94	0.50
39:DQ:60:LEU:HD23	39:DQ:60:LEU:C	2.32	0.50
23:DA:2050:C:H1'	26:DD:156:MET:HE1	1.93	0.50
23:BA:1827:C:H2'	23:BA:1828:G:O4'	2.12	0.49
25:BC:132:PRO:HG3	25:BC:190:TYR:CE1	2.46	0.49
44:BV:76:LEU:CD1	44:BV:76:LEU:H	2.24	0.49
10:AJ:82:ILE:O	10:AJ:86:MET:HB2	2.12	0.49
1:CA:80:G:H2'	1:CA:81:G:H8	1.77	0.49
26:DD:9:VAL:HG13	26:DD:25:VAL:O	2.11	0.49
23:DA:661:C:H4'	34:DL:16:ARG:CD	2.42	0.49
3:AC:6:HIS:CD2	3:AC:7:PRO:HD2	2.43	0.49
30:DH:77:LEU:O	30:DH:143:SER:HB3	2.12	0.49
23:BA:774:A:H2	23:BA:787:U:O2'	1.93	0.49
46:DX:90:ILE:O	46:DX:94:LEU:HB2	2.12	0.49
33:BK:2:ILE:CD1	33:BK:82:ASN:HD22	2.25	0.49
23:BA:2894:G:N3	23:BA:2894:G:H2'	2.27	0.49
1:CA:1520:G:H2'	1:CA:1521:G:H8	1.77	0.49
1:CA:114:U:H2'	1:CA:115:G:H8	1.77	0.49
23:DA:1486:A:N6	23:DA:1504:C:H42	2.10	0.49
23:BA:581:C:H2'	23:BA:582:G:C8	2.47	0.49
17:AQ:59:ILE:HG22	17:AQ:60:ILE:N	2.27	0.49
25:DC:72:LYS:CE	25:DC:101:GLU:HG2	2.42	0.49
17:CQ:21:VAL:HG11	17:CQ:59:ILE:HD11	1.93	0.49
23:BA:903:C:H2'	23:BA:904:C:C6	2.47	0.49
1:CA:345:C:OP2	38:DP:39:ARG:NH2	2.45	0.49
43:BU:29:GLU:HB3	43:BU:38:ILE:HB	1.94	0.49
23:DA:2208:U:O2'	23:DA:2209:C:H5'	2.12	0.49
23:BA:496:G:H1'	41:BS:61:ASN:ND2	2.27	0.49
31:DI:9:LEU:O	31:DI:13:LEU:HG	2.12	0.49
16:AP:23:ASP:O	16:AP:26:ARG:HB2	2.11	0.49
47:BY:11:GLU:N	47:BY:11:GLU:OE1	2.43	0.49
23:DA:1111:A:N3	23:DA:1112:G:H1'	2.27	0.49
23:BA:1165:U:H2'	23:BA:1166:C:C6	2.47	0.49
39:DQ:79:PHE:CD1	39:DQ:79:PHE:C	2.86	0.49
23:DA:1161:C:O2'	40:DR:23:GLU:HG2	2.11	0.49
39:BQ:79:PHE:C	39:BQ:79:PHE:CD1	2.85	0.49
39:BQ:88:ILE:HB	39:BQ:90:VAL:CG1	2.36	0.49
23:DA:114(B):A:N3	23:DA:1144:G:C8	2.80	0.49
23:BA:2246:G:H2'	23:BA:2247:A:C8	2.48	0.49
23:DA:2392:A:H2	23:DA:2424:C:N4	2.08	0.49
3:AC:105:GLU:HG2	3:AC:106:VAL:N	2.18	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BA:1578:U:H2'	23:BA:1579:A:H5''	1.94	0.49
3:CC:14:ILE:HG23	3:CC:15:THR:N	2.25	0.49
2:CB:187:LEU:HD11	2:CB:204:ASN:O	2.12	0.49
26:DD:51:PHE:HB3	26:DD:52:LEU:HD12	1.94	0.49
25:DC:244:ARG:HB2	25:DC:244:ARG:HH11	1.77	0.49
30:DH:82:ARG:HG2	30:DH:89:TYR:CD1	2.47	0.49
24:DB:71:C:C4	24:DB:72:G:N7	2.80	0.49
1:CA:1152:A:H2'	1:CA:1153:C:C6	2.46	0.49
1:AA:1295:G:O2'	13:AM:14:ARG:CZ	2.60	0.49
23:BA:390:A:C6	34:BL:71:VAL:HG21	2.47	0.49
4:AD:122:ARG:HD3	4:AD:122:ARG:O	2.12	0.49
26:DD:120:TRP:CD1	26:DD:155:LYS:HB3	2.47	0.49
19:CS:63:THR:HG22	19:CS:66:MET:HE3	1.94	0.49
15:AO:39:LEU:HD12	15:AO:56:LEU:HB2	1.93	0.49
1:CA:724:G:C2	1:CA:725:G:C8	3.00	0.49
1:CA:971:G:C8	1:CA:1365:G:H4'	2.48	0.49
44:BV:82:ARG:HG2	44:BV:83:PRO:HD2	1.94	0.49
33:DK:24:VAL:CG2	33:DK:33:ALA:HB2	2.41	0.49
23:BA:144:C:H2'	23:BA:145:G:H8	1.77	0.49
40:DR:14:VAL:HG13	40:DR:96:ILE:HG13	1.94	0.49
2:CB:61:LEU:HD21	2:CB:68:ILE:HG12	1.95	0.49
23:BA:2419:U:O4	53:B5:30:ARG:CZ	2.60	0.49
23:DA:270(Q):C:HO2'	23:DA:270(R):C:H6	1.58	0.49
23:DA:2853:C:H2'	23:DA:2854:G:H8	1.77	0.49
28:BF:16:ARG:O	28:BF:20:ILE:HG12	2.12	0.49
1:CA:1203:C:H2'	1:CA:1204:A:H8	1.77	0.49
28:DF:153:ARG:HB3	28:DF:153:ARG:NH1	2.27	0.49
33:DK:61:VAL:O	33:DK:61:VAL:HG13	2.11	0.49
23:BA:1689:A:H62	23:BA:1698:A:H2	1.59	0.49
23:BA:114(B):A:N3	23:BA:1144:G:C8	2.81	0.49
23:DA:827:U:O2	23:DA:2246:G:H4'	2.12	0.49
23:BA:2394:C:H2'	23:BA:2395:C:C6	2.48	0.49
23:DA:114(B):A:C4	23:DA:1144:G:N7	2.80	0.49
1:AA:1286:A:N7	21:AU:22:ARG:NH2	2.61	0.49
23:BA:956:G:N2	23:BA:959:A:H3'	2.26	0.49
46:BX:11:ARG:HB2	46:BX:13:ILE:HG22	1.94	0.49
13:AM:98:VAL:HB	13:AM:99:ARG:HH11	1.77	0.49
23:BA:603:A:N6	23:BA:655:A:H1'	2.27	0.49
23:DA:603:A:H61	23:DA:655:A:C1'	2.25	0.49
23:BA:1568:G:OP2	25:BC:63:ARG:NH2	2.45	0.49
23:DA:1431:U:H2'	23:DA:1432:C:C6	2.47	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:AC:175:LEU:O	3:AC:175:LEU:HD23	2.11	0.49
35:DM:60:ARG:N	44:DV:179:ASP:HB2	2.27	0.49
23:BA:860:U:O4'	23:BA:860:U:O2	2.29	0.49
43:BU:20:TYR:CE1	43:BU:42:VAL:HA	2.47	0.49
1:AA:1226:C:H2'	13:AM:103:THR:HB	1.93	0.49
23:DA:1478:G:H2'	23:DA:1479:G:H8	1.77	0.49
25:BC:10:THR:HG23	25:BC:13:ARG:HB3	1.94	0.49
30:BH:57:ARG:O	30:BH:61:ARG:HG3	2.12	0.49
1:CA:1157:A:N6	1:CA:1178:G:H1'	2.27	0.49
36:BN:103:ARG:HH12	36:BN:110:PRO:HG3	1.78	0.49
23:DA:2335:A:H2'	37:DO:13:ARG:HH22	1.77	0.49
1:AA:1290:G:H2'	1:AA:1290:G:N3	2.25	0.49
23:BA:2010:G:H5''	41:BS:42:ARG:HB2	1.94	0.49
27:BE:143:ALA:HB1	27:BE:148:LEU:HB2	1.94	0.49
23:DA:165:U:H2'	23:DA:171:G:O4'	2.13	0.49
1:CA:665:A:H2'	1:CA:732:C:O2	2.12	0.49
23:DA:903:C:H2'	23:DA:904:C:C6	2.47	0.49
24:BB:63:G:H2'	24:BB:64:C:C6	2.47	0.49
16:AP:8:ARG:HB3	16:AP:28:ARG:NH1	2.26	0.49
33:BK:112:MET:HA	33:BK:115:VAL:HG22	1.93	0.49
2:AB:145:LEU:O	2:AB:149:LEU:HB2	2.12	0.49
2:AB:75:LYS:C	2:AB:75:LYS:HD3	2.32	0.49
23:BA:150:C:H2'	23:BA:151:C:C6	2.47	0.49
23:DA:1971:A:C4	25:DC:241:PRO:HG3	2.47	0.49
23:BA:114(B):A:C4	23:BA:1144:G:N7	2.80	0.49
53:B5:33:ASN:ND2	53:B5:34:TRP:H	2.11	0.49
19:CS:16:LEU:HA	19:CS:19:VAL:HG12	1.94	0.49
10:CJ:33:GLN:O	10:CJ:75:ILE:HG12	2.12	0.49
53:D5:14:VAL:CG1	53:D5:22:VAL:HG13	2.42	0.49
25:BC:81:ALA:HB3	25:BC:94:LEU:HB3	1.95	0.49
37:DO:14:VAL:O	37:DO:18:ILE:HG12	2.13	0.49
23:DA:2516:G:C6	23:DA:2517:C:N4	2.80	0.49
26:DD:117:MET:CE	26:DD:136:ARG:HA	2.41	0.49
23:DA:2481:G:HO2'	23:DA:2482:G:P	2.35	0.49
27:DE:155:LEU:CD2	27:DE:186:ILE:HD13	2.40	0.49
29:DG:51:ARG:O	29:DG:52:VAL:HG23	2.13	0.49
35:BM:60:ARG:N	44:BV:179:ASP:HB2	2.26	0.49
51:D3:36:LEU:HB3	51:D3:50:ARG:NH1	2.27	0.49
1:AA:1014:A:H4'	19:AS:14:HIS:CE1	2.46	0.49
22:CV:6171:U:H2'	22:CV:6172:U:C6	2.47	0.49
24:DB:75:G:N1	24:DB:102:G:N2	2.60	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:DC:10:THR:HG23	25:DC:13:ARG:HB2	1.94	0.49
23:BA:1937:A:N7	23:BA:1939:U:H2'	2.28	0.49
33:BK:2:ILE:CG1	33:BK:8:LEU:HD11	2.42	0.49
23:DA:2894:G:H2'	23:DA:2894:G:N3	2.27	0.49
23:DA:966:G:C4	23:DA:967:C:C5	3.00	0.49
23:DA:1486:A:C6	23:DA:1504:C:N4	2.80	0.49
51:B3:13:CYS:O	51:B3:21:TYR:HA	2.13	0.49
41:DS:12:ILE:HD12	41:DS:46:PHE:CE2	2.48	0.49
35:BM:52:VAL:O	35:BM:56:ARG:HB2	2.12	0.49
1:CA:1305:G:H1'	1:CA:1306:A:C8	2.48	0.49
51:D3:16:CYS:SG	51:D3:48:VAL:HG23	2.53	0.49
9:CI:4:TYR:CE2	9:CI:88:TYR:HB2	2.47	0.49
38:DP:24:PRO:HA	38:DP:49:VAL:HG13	1.93	0.49
35:BM:23:GLY:HA3	35:BM:98:LYS:HB2	1.94	0.49
32:DJ:77:VAL:HB	32:DJ:145:VAL:HG22	1.93	0.49
1:CA:444:C:H2'	1:CA:445:G:C8	2.47	0.49
8:CH:1:MET:HE2	8:CH:1:MET:N	2.28	0.49
1:AA:767:A:H2'	1:AA:768:A:O4'	2.12	0.49
23:BA:754:C:H2'	23:BA:755:C:C6	2.47	0.49
23:DA:41:C:H2'	23:DA:43:G:O4'	2.11	0.49
1:CA:1391:U:H2'	1:CA:1392:G:C8	2.47	0.49
23:DA:471:A:H2'	23:DA:472:A:O4'	2.12	0.49
23:DA:855:G:H2'	23:DA:856:C:C6	2.47	0.49
51:B3:16:CYS:SG	51:B3:48:VAL:HG23	2.52	0.49
9:AI:19:LEU:HD23	9:AI:20:ARG:N	2.26	0.49
23:BA:1826:G:OP1	25:BC:233:HIS:HD2	1.96	0.49
1:CA:89:U:H2'	1:CA:90:C:C6	2.47	0.49
35:BM:76:LYS:H	35:BM:88:GLY:HA2	1.76	0.49
23:DA:2439:A:H8	23:DA:2439:A:H5'	1.70	0.49
29:DG:92:ILE:H	29:DG:92:ILE:CD1	2.21	0.49
2:CB:20:GLU:HA	2:CB:20:GLU:OE1	2.12	0.49
2:CB:24:TRP:HE3	2:CB:25:ASN:O	1.96	0.49
26:DD:201:THR:O	26:DD:202:LYS:HD3	2.13	0.49
28:BF:86:MET:H	28:BF:87:PRO:CD	2.25	0.49
11:AK:20:TYR:O	11:AK:30:VAL:HA	2.13	0.49
51:B3:36:LEU:HB3	51:B3:50:ARG:NH1	2.27	0.49
25:BC:79:VAL:O	25:BC:113:VAL:HG13	2.11	0.49
23:DA:1478:G:HO2'	23:DA:1558:A:H2	1.61	0.49
2:CB:8:LYS:HG2	2:CB:217:ARG:NH1	2.27	0.49
23:DA:596:G:C6	23:DA:597:U:C4	3.01	0.49
43:DU:90:LEU:HG	43:DU:91:GLU:H	1.76	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:626:U:H2'	1:AA:627:G:H8	1.76	0.49
23:BA:1437:C:H2'	23:BA:1438:U:H6	1.78	0.49
27:BE:125:LEU:HB3	27:BE:196:LEU:HD23	1.94	0.49
23:BA:2564:A:OP1	23:BA:2648:C:H4'	2.13	0.49
36:BN:87:TYR:OH	36:BN:116:LEU:HB3	2.13	0.49
10:CJ:30:SER:HB2	10:CJ:80:LYS:HG2	1.93	0.49
23:BA:23:G:H2'	23:BA:24:G:H8	1.77	0.49
1:CA:1366:C:H2'	1:CA:1367:C:C6	2.48	0.49
23:DA:1354:A:C8	23:DA:1355:G:C8	3.01	0.49
4:AD:195:ALA:C	4:AD:196:LEU:HD12	2.32	0.49
28:BF:133:LEU:HD21	28:BF:157:ILE:HG13	1.94	0.49
4:AD:43:HIS:HA	4:AD:46:LYS:HE3	1.94	0.49
12:CL:57:VAL:O	12:CL:59:LEU:HD22	2.13	0.49
1:AA:1413:A:C6	1:AA:1414:U:C4	3.00	0.49
23:BA:1592:C:H2'	23:BA:1593:G:H8	1.78	0.49
23:BA:405:U:H3'	23:BA:406:G:H5'	1.94	0.49
23:DA:1963:U:O2	23:DA:1963:U:H2'	2.13	0.49
2:AB:22:LYS:HZ3	2:AB:22:LYS:H	1.60	0.49
10:CJ:29:ARG:HG2	10:CJ:29:ARG:O	2.12	0.49
4:CD:53:ASP:O	4:CD:57:ARG:HD3	2.13	0.49
11:AK:17:GLY:HA3	11:AK:77:MET:SD	2.52	0.49
23:DA:2340:G:H2'	23:DA:2341:G:H8	1.77	0.49
30:BH:92:VAL:HG22	30:BH:120:ILE:HD12	1.94	0.49
23:BA:1970:A:H4'	23:BA:1971:A:OP1	2.12	0.49
23:BA:1022:G:C6	23:BA:1140:C:C4	3.01	0.49
23:BA:733:G:H8	23:BA:733:G:O5'	1.96	0.49
1:CA:979:C:N4	14:CN:18:VAL:HG12	2.28	0.49
50:D2:33:CYS:SG	50:D2:49:CYS:SG	3.07	0.49
12:CL:76:LEU:HD11	12:CL:106:ALA:HA	1.95	0.49
3:CC:17:ASP:CB	3:CC:21:ARG:HH22	2.23	0.49
26:BD:5:LEU:HB2	26:BD:51:PHE:HD2	1.77	0.49
23:BA:557:U:H2'	23:BA:558:G:H8	1.76	0.49
23:DA:322:A:OP2	27:DE:169:ASN:HB2	2.12	0.49
23:DA:1188:U:C2'	23:DA:1189:A:H5'	2.42	0.49
1:CA:1080:A:H5''	1:CA:1081:G:OP2	2.12	0.49
23:DA:1608:A:HO2'	23:DA:1610:A:P	2.35	0.49
23:BA:596:G:C6	23:BA:597:U:C4	3.01	0.49
25:BC:108:PRO:CG	25:BC:143:HIS:CE1	2.96	0.49
36:DN:81:ASP:O	36:DN:85:PRO:HG2	2.12	0.49
34:DL:126:VAL:HA	34:DL:145:PRO:HB2	1.95	0.49
28:BF:130:ASN:OD1	28:BF:160:VAL:HA	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
32:BJ:80:ALA:HB3	32:BJ:147:ALA:HB2	1.95	0.49
11:CK:124:LYS:C	11:CK:126:ARG:H	2.16	0.49
1:CA:624:C:H2'	1:CA:625:G:H8	1.78	0.49
23:BA:1833:U:C2	23:BA:1834:U:C6	3.00	0.49
17:AQ:59:ILE:CG2	17:AQ:71:PHE:HB3	2.42	0.49
20:AT:85:MET:HB2	20:AT:104:LEU:HD21	1.94	0.49
23:DA:144:C:H2'	23:DA:145:G:H8	1.77	0.49
18:AR:43:PHE:O	18:AR:51:LEU:HB2	2.12	0.49
29:BG:23:ARG:N	29:BG:23:ARG:HD3	2.27	0.49
23:DA:118:A:N3	23:DA:178:G:H1'	2.27	0.49
19:CS:53:ASN:HD21	19:CS:56:GLN:H	1.59	0.49
23:BA:2258:C:H4'	23:BA:2259:G:OP2	2.12	0.49
3:CC:79:ARG:N	3:CC:79:ARG:HD3	2.27	0.49
1:CA:488:C:O5'	1:CA:488:C:H6	1.94	0.49
23:DA:69:C:H2'	23:DA:70:G:C8	2.47	0.49
34:DL:32:THR:OG1	34:DL:36:LYS:HB3	2.13	0.49
45:DW:37:LEU:HG	45:DW:60:PHE:HA	1.95	0.49
47:BY:14:ARG:HA	47:BY:17:SER:HB2	1.94	0.49
10:CJ:82:ILE:O	10:CJ:86:MET:HB2	2.12	0.49
1:CA:134:A:H61	16:CP:25:ARG:HH12	1.56	0.49
23:DA:603:A:N6	23:DA:655:A:H1'	2.28	0.49
28:DF:98:ARG:O	28:DF:101:ILE:HG12	2.12	0.49
11:AK:21:ILE:HB	11:AK:84:VAL:HG12	1.94	0.49
35:BM:43:THR:OG1	35:BM:46:GLN:HG3	2.12	0.49
23:BA:795:C:O2'	23:BA:796:C:H5'	2.13	0.49
3:CC:19:GLU:HG3	3:CC:54:ARG:HD2	1.95	0.49
1:CA:1349:A:H2'	1:CA:1350:A:H8	1.76	0.49
1:AA:475:G:H2'	1:AA:476:G:C8	2.45	0.49
23:BA:2892:A:N6	23:BA:2893:G:C2	2.81	0.49
7:AG:41:ARG:NH1	7:AG:41:ARG:HB3	2.28	0.49
20:AT:49:ALA:HB3	20:AT:99:LEU:HD12	1.94	0.49
2:CB:17:PHE:HB2	2:CB:42:ILE:HG22	1.95	0.49
17:AQ:57:VAL:HG12	17:AQ:76:LEU:HA	1.95	0.49
1:AA:909:A:H2'	1:AA:910:C:O4'	2.13	0.49
33:DK:88:ASN:OD1	33:DK:89:ASN:N	2.46	0.49
33:BK:12:ASP:OD1	33:BK:85:VAL:HG13	2.13	0.49
45:BW:50:ASN:HD22	45:BW:83:PRO:HD3	1.78	0.49
23:DA:1887:C:H3'	23:DA:1888:G:H5''	1.95	0.49
23:DA:2784:C:H2'	23:DA:2785:C:C6	2.47	0.49
4:CD:88:VAL:O	4:CD:92:VAL:HG23	2.12	0.49
23:DA:258:G:H2'	23:DA:259:G:H8	1.78	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:AD:162:LEU:HD13	4:AD:181:MET:HG2	1.95	0.49
1:AA:1300:G:O2'	1:AA:1301:U:P	2.71	0.49
21:CU:6:ARG:NE	21:CU:15:ARG:HH12	2.10	0.49
23:DA:289:A:H2'	23:DA:290:G:O4'	2.13	0.49
23:BA:1973:G:H2'	23:BA:1974:C:H6	1.77	0.49
6:CF:26:ILE:O	6:CF:30:LEU:HG	2.13	0.49
23:DA:116:C:H2'	23:DA:117:G:C8	2.47	0.49
1:AA:19:C:H5''	5:AE:86:ALA:HB1	1.95	0.49
45:BW:53:MET:HA	45:BW:58:THR:O	2.13	0.49
3:CC:113:ALA:HB3	3:CC:114:PRO:HD3	1.95	0.49
23:DA:405:U:H3'	23:DA:406:G:H5'	1.95	0.49
23:DA:943:U:OP2	34:DL:38:GLN:CD	2.51	0.49
34:DL:35:HIS:O	34:DL:36:LYS:HB2	2.13	0.49
1:CA:1443:G:H22	38:DP:119:LYS:HA	1.78	0.49
23:DA:1022:G:O2'	23:DA:1023:U:OP2	2.28	0.49
23:BA:783:A:C8	23:BA:783:A:H3'	2.48	0.49
30:BH:5:LEU:CD2	30:BH:5:LEU:H	2.18	0.49
10:AJ:32:ALA:CB	10:AJ:76:ASN:HB2	2.43	0.49
3:AC:195:VAL:CG1	3:AC:196:LEU:H	2.20	0.49
26:BD:117:MET:CE	26:BD:136:ARG:HA	2.38	0.49
2:CB:28:PHE:CD1	2:CB:190:THR:HA	2.48	0.49
26:DD:169:ASN:ND2	26:DD:201:THR:HG21	2.28	0.49
19:AS:32:LYS:HA	19:AS:50:ALA:HB3	1.95	0.49
2:CB:88:ALA:HA	2:CB:223:ILE:HD11	1.95	0.49
23:DA:2365:G:O6	53:D5:39:LYS:HE3	2.13	0.49
23:BA:1486:A:C6	23:BA:1504:C:N4	2.80	0.49
25:DC:10:THR:HG23	25:DC:13:ARG:HB3	1.93	0.49
51:D3:11:LEU:HD21	51:D3:51:GLU:CD	2.33	0.49
1:CA:255:G:H2'	1:CA:256:U:H6	1.78	0.49
2:AB:8:LYS:HG2	2:AB:217:ARG:NH1	2.28	0.49
1:CA:1503:A:H5''	1:CA:1531:A:H1'	1.94	0.49
52:B4:21:ARG:HB3	52:B4:31:LEU:CD2	2.43	0.49
51:D3:13:CYS:O	51:D3:21:TYR:HA	2.13	0.49
1:AA:1157:A:N6	1:AA:1178:G:H1'	2.27	0.49
3:AC:73:PRO:O	3:AC:76:VAL:HG22	2.13	0.49
44:DV:82:ARG:HG2	44:DV:83:PRO:HD2	1.95	0.49
20:CT:49:ALA:HB3	20:CT:99:LEU:HD12	1.94	0.49
1:CA:1154:G:H2'	1:CA:1155:G:C8	2.48	0.49
8:AH:110:ALA:HB3	8:AH:121:ASP:HB3	1.94	0.49
45:DW:53:MET:HA	45:DW:58:THR:O	2.12	0.49
19:AS:51:VAL:O	19:AS:58:VAL:HG22	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BA:2537:U:H2'	23:BA:2538:C:H6	1.77	0.49
2:AB:212:GLN:HE22	2:AB:216:SER:HB2	1.77	0.49
25:BC:105:ILE:CG1	25:BC:106:ILE:HD12	2.42	0.49
4:CD:43:HIS:HA	4:CD:46:LYS:HE3	1.93	0.49
23:BA:2012:G:O2'	41:BS:96:ILE:HD11	2.12	0.49
28:BF:153:ARG:HB3	28:BF:153:ARG:NH1	2.27	0.49
23:DA:1027:A:C2	23:DA:2488:A:H5'	2.47	0.49
2:CB:212:GLN:HE22	2:CB:216:SER:HB2	1.78	0.49
23:DA:2512:C:H5''	23:DA:2513:G:OP2	2.13	0.49
1:CA:538:G:O3'	12:CL:113:LYS:HG3	2.13	0.49
23:BA:747:U:O2	23:BA:2014:A:H1'	2.12	0.49
46:BX:13:ILE:HG23	46:BX:14:VAL:H	1.77	0.49
36:DN:50:HIS:O	36:DN:54:LEU:HB2	2.13	0.49
23:DA:1171:G:H2'	23:DA:1173:G:O4'	2.13	0.49
26:BD:101:ARG:HG2	26:BD:171:GLU:HA	1.95	0.49
3:AC:14:ILE:HG23	3:AC:15:THR:N	2.25	0.49
28:DF:86:MET:H	28:DF:87:PRO:CD	2.26	0.49
1:AA:1313:U:OP1	19:AS:6:LYS:HG3	2.12	0.49
23:DA:1517:G:H2'	23:DA:1518:C:C6	2.48	0.49
3:AC:6:HIS:ND1	14:AN:49:HIS:HB3	2.28	0.49
23:BA:1614:A:C6	41:BS:87:PRO:HA	2.48	0.49
1:CA:1194:U:H4'	5:CE:22:GLY:O	2.12	0.49
19:CS:5:LEU:HG	19:CS:10:PHE:HB3	1.95	0.49
46:BX:58:ILE:HD11	46:BX:91:LYS:CG	2.43	0.49
2:AB:118:LEU:O	2:AB:122:PHE:HB2	2.13	0.49
1:AA:724:G:C2	1:AA:725:G:C8	3.00	0.49
23:DA:1495:A:H2'	23:DA:1496:A:C2	2.48	0.49
17:AQ:76:LEU:HD12	17:AQ:77:VAL:H	1.77	0.49
1:AA:1100:C:OP2	2:AB:96:ARG:HG2	2.13	0.49
30:DH:115:ALA:HB3	30:DH:129:THR:O	2.12	0.49
23:BA:110:G:C2	23:BA:111:A:C8	3.01	0.49
19:CS:51:VAL:O	19:CS:58:VAL:HG22	2.12	0.49
23:DA:466:A:N3	23:DA:683:C:H1'	2.28	0.49
23:BA:1963:U:H2'	23:BA:1963:U:O2	2.12	0.49
17:AQ:52:LYS:HD2	17:AQ:52:LYS:H	1.77	0.49
1:CA:509:A:C6	1:CA:510:A:N1	2.81	0.49
23:BA:433:C:H2'	23:BA:434:U:C6	2.48	0.49
39:DQ:79:PHE:HD1	39:DQ:79:PHE:C	2.16	0.49
40:DR:47:VAL:HG13	40:DR:52:VAL:N	2.27	0.49
40:BR:47:VAL:HG12	40:BR:49:THR:O	2.13	0.49
25:DC:242:ARG:CD	25:DC:242:ARG:N	2.71	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DM:76:LYS:H	35:DM:88:GLY:HA2	1.76	0.49
11:CK:44:SER:OG	11:CK:47:VAL:HG23	2.12	0.49
25:DC:31:LYS:HE3	25:DC:33:LEU:HD21	1.94	0.49
25:BC:25:THR:HG21	25:BC:81:ALA:CA	2.43	0.49
26:DD:201:THR:CG2	26:DD:202:LYS:H	2.22	0.49
12:CL:82:VAL:HG21	12:CL:99:ILE:CD1	2.43	0.49
32:DJ:66:THR:N	32:DJ:71:MET:HE3	2.22	0.49
6:CF:19:LEU:O	6:CF:23:LYS:HG3	2.13	0.49
4:CD:31:CYS:O	4:CD:32:ALA:HB3	2.12	0.49
3:AC:19:GLU:HG3	3:AC:54:ARG:HD2	1.95	0.49
23:BA:1486:A:N6	23:BA:1504:C:H42	2.09	0.49
4:AD:163:GLU:O	4:AD:166:LYS:HG3	2.13	0.49
25:DC:70:TRP:CH2	25:DC:150:LYS:HA	2.48	0.49
9:AI:99:LEU:HD12	9:AI:101:PHE:HE2	1.77	0.49
34:BL:95:VAL:HG23	34:BL:125:VAL:HA	1.94	0.49
13:AM:49:THR:O	13:AM:53:VAL:HG23	2.12	0.49
33:DK:79:PHE:HD2	38:DP:72:VAL:HG22	1.78	0.49
7:CG:41:ARG:NH1	7:CG:41:ARG:HB3	2.28	0.49
5:AE:12:LEU:HD22	5:AE:12:LEU:C	2.32	0.49
25:BC:72:LYS:HE3	25:BC:101:GLU:CB	2.43	0.49
23:DA:69:C:O2'	23:DA:70:G:H5'	2.13	0.49
38:BP:107:ASP:O	38:BP:110:ILE:HG22	2.12	0.49
11:CK:17:GLY:HA3	11:CK:77:MET:SD	2.52	0.49
23:BA:2784:C:H2'	23:BA:2785:C:C6	2.48	0.49
23:DA:634:C:H2'	23:DA:635:C:C6	2.48	0.49
1:CA:565:U:C6	1:CA:566:G:C8	3.01	0.49
23:BA:1680:U:O2	23:BA:1763:G:H3'	2.12	0.49
32:DJ:58:ARG:C	32:DJ:60:LYS:H	2.15	0.49
1:AA:865:A:O5'	1:AA:865:A:H8	1.96	0.49
26:DD:14:ILE:HD12	26:DD:14:ILE:C	2.33	0.49
32:BJ:160:LYS:CD	32:BJ:161:LEU:H	2.26	0.49
50:D2:3:LYS:N	50:D2:3:LYS:HD2	2.28	0.48
50:B2:3:LYS:O	50:B2:4:HIS:C	2.51	0.48
39:BQ:79:PHE:C	39:BQ:79:PHE:HD1	2.16	0.48
23:DA:733:G:O5'	23:DA:733:G:H8	1.96	0.48
43:BU:14:LEU:HD23	43:BU:15:VAL:N	2.28	0.48
43:DU:2:ARG:C	43:DU:4:LYS:H	2.17	0.48
45:DW:32:ARG:C	45:DW:35:ASN:HD21	2.15	0.48
28:BF:55:LYS:HD2	28:BF:58:GLN:NE2	2.20	0.48
23:BA:947:G:N2	23:BA:971:C:C2	2.81	0.48
16:AP:27:LYS:N	16:AP:27:LYS:HD2	2.26	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BA:1431:U:H2'	23:BA:1432:C:C6	2.48	0.48
23:BA:2741:A:H2'	23:BA:2742:C:O4'	2.13	0.48
28:DF:173:LEU:HA	28:DF:176:LEU:HD12	1.93	0.48
3:CC:175:LEU:O	3:CC:175:LEU:HD23	2.12	0.48
18:CR:50:ILE:HD11	18:CR:74:ARG:NH1	2.28	0.48
51:B3:11:LEU:HD21	51:B3:51:GLU:CD	2.33	0.48
23:BA:322:A:H3'	27:BE:169:ASN:HD21	1.78	0.48
1:CA:427:U:C4	1:CA:428:G:C6	3.00	0.48
1:AA:427:U:C4	1:AA:428:G:C6	3.01	0.48
23:BA:481:G:O2'	23:BA:482:A:P	2.71	0.48
20:AT:67:ALA:HA	20:AT:72:LEU:O	2.12	0.48
23:DA:1437:C:H2'	23:DA:1438:U:H6	1.75	0.48
1:CA:600:C:H2'	1:CA:601:C:C6	2.48	0.48
18:CR:56:THR:HB	18:CR:58:LEU:HD13	1.94	0.48
23:DA:784:A:N7	25:DC:229:VAL:HG21	2.27	0.48
38:BP:54:ARG:HA	38:BP:59:THR:OG1	2.12	0.48
27:DE:12:LEU:HB2	27:DE:124:LEU:HD11	1.94	0.48
3:CC:73:PRO:O	3:CC:76:VAL:HG22	2.13	0.48
1:CA:1237:C:OP1	1:CA:1238:A:H1'	2.13	0.48
25:BC:72:LYS:CE	25:BC:101:GLU:HG2	2.42	0.48
17:CQ:59:ILE:HG22	17:CQ:60:ILE:N	2.28	0.48
23:DA:2795:G:H3'	23:DA:2797:U:H5''	1.95	0.48
23:DA:1615:C:O2'	23:DA:1616:A:H5'	2.13	0.48
23:BA:1270:C:H5''	23:BA:1271:G:O5'	2.13	0.48
1:CA:1300:G:O2'	1:CA:1301:U:P	2.71	0.48
23:BA:1001:A:H2'	23:BA:1002:G:O4'	2.12	0.48
8:AH:81:HIS:HB2	8:AH:138:TRP:OXT	2.11	0.48
12:AL:57:VAL:O	12:AL:59:LEU:HD22	2.12	0.48
23:DA:329:G:H1	43:DU:19:LYS:HE3	1.78	0.48
23:BA:2252:G:H2'	23:BA:2253:G:H8	1.78	0.48
47:BY:1:MET:O	47:BY:1:MET:SD	2.71	0.48
23:BA:165:U:H2'	23:BA:171:G:O4'	2.13	0.48
23:DA:2436:G:C5	23:DA:2437:U:C5	3.01	0.48
40:BR:47:VAL:HG13	40:BR:52:VAL:N	2.28	0.48
23:DA:114(B):A:C4	23:DA:1144:G:C8	3.01	0.48
19:AS:16:LEU:HA	19:AS:19:VAL:HG12	1.94	0.48
44:DV:99:TYR:CE2	44:DV:125:LEU:HD12	2.48	0.48
10:CJ:32:ALA:CB	10:CJ:76:ASN:HB2	2.43	0.48
13:CM:98:VAL:HB	13:CM:99:ARG:HH11	1.78	0.48
1:AA:80:G:H2'	1:AA:81:G:H8	1.78	0.48
1:AA:1378:C:H5''	7:AG:6:ARG:NE	2.25	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BA:2039:C:C2	23:BA:2040:C:C5	3.01	0.48
30:BH:77:LEU:O	30:BH:143:SER:HB3	2.13	0.48
35:DM:43:THR:OG1	35:DM:46:GLN:HG3	2.13	0.48
23:BA:2415:G:H4'	34:BL:66:GLY:HA2	1.92	0.48
1:CA:277:C:OP1	17:CQ:41:LYS:HE3	2.12	0.48
23:DA:1937:A:N7	23:DA:1939:U:H2'	2.27	0.48
35:BM:39:PRO:O	35:BM:40:ALA:HB2	2.13	0.48
25:BC:43:ARG:HB2	25:BC:49:ILE:HA	1.96	0.48
38:BP:58:ASN:C	38:BP:58:ASN:HD22	2.17	0.48
11:AK:105:VAL:HG23	11:AK:105:VAL:O	2.13	0.48
1:CA:1412:C:H2'	1:CA:1413:A:H8	1.77	0.48
23:DA:380:U:O2	46:DX:20:ARG:NH2	2.46	0.48
1:AA:1412:C:H2'	1:AA:1413:A:C8	2.48	0.48
34:BL:138:LEU:HD11	34:BL:144:GLU:HB3	1.94	0.48
1:AA:112:G:C2	1:AA:113:G:C8	3.02	0.48
1:CA:832:C:N4	1:CA:855:G:O6	2.47	0.48
23:DA:718:A:O5'	23:DA:718:A:H8	1.96	0.48
23:BA:1252:G:C2	23:BA:1253:A:C2	3.01	0.48
23:BA:996:A:H4'	39:BQ:92:ARG:CZ	2.44	0.48
34:DL:114:ILE:CD1	34:DL:130:PHE:CD1	2.90	0.48
12:AL:26:LEU:HD13	12:AL:27:LYS:N	2.17	0.48
23:DA:270(H):C:H2'	23:DA:270(I):C:H6	1.77	0.48
28:DF:55:LYS:O	28:DF:59:GLU:HG3	2.13	0.48
10:CJ:33:GLN:HB2	10:CJ:75:ILE:HD11	1.95	0.48
23:DA:124:G:N7	52:D4:19:ARG:NH2	2.61	0.48
23:BA:1171:G:H2'	23:BA:1173:G:O4'	2.13	0.48
23:DA:2516:G:O6	23:DA:2517:C:N4	2.47	0.48
19:AS:34:TRP:CZ2	19:AS:57:HIS:HE1	2.31	0.48
10:AJ:34:VAL:CG2	10:AJ:74:ILE:HG22	2.43	0.48
30:DH:88:ILE:CG2	30:DH:89:TYR:N	2.76	0.48
37:BO:25:ARG:HD2	37:BO:88:ASP:OD1	2.12	0.48
3:CC:6:HIS:ND1	14:CN:49:HIS:HB3	2.28	0.48
43:DU:76:CYS:CB	43:DU:77:PRO:CD	2.91	0.48
43:BU:90:LEU:HG	43:BU:91:GLU:H	1.77	0.48
23:DA:1953:A:C2	23:DA:2549:G:N3	2.79	0.48
1:CA:1062:U:H2'	1:CA:1063:C:C6	2.49	0.48
1:AA:1503:A:H5''	1:AA:1531:A:H1'	1.93	0.48
23:DA:1567:A:H5''	25:DC:58:HIS:CD2	2.48	0.48
1:AA:600:C:H2'	1:AA:601:C:C6	2.48	0.48
23:BA:2699:C:H2'	23:BA:2700:C:O4'	2.13	0.48
1:AA:1237:C:OP1	1:AA:1238:A:H1'	2.12	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:971:G:C8	1:AA:1365:G:H4'	2.48	0.48
44:DV:10:ARG:HH21	44:DV:26:GLY:H	1.62	0.48
5:AE:70:PRO:HB3	5:AE:144:THR:HG22	1.95	0.48
26:DD:31:CYS:HB3	26:DD:49:LEU:HB3	1.95	0.48
4:AD:53:ASP:O	4:AD:57:ARG:HD3	2.14	0.48
23:DA:1001:A:H2'	23:DA:1002:G:O4'	2.13	0.48
23:BA:1632:A:O5'	23:BA:1632:A:H8	1.96	0.48
12:CL:19:LYS:H	12:CL:19:LYS:HD3	1.78	0.48
1:CA:45:U:O5'	1:CA:45:U:H6	1.97	0.48
23:DA:496:G:H1'	41:DS:61:ASN:ND2	2.29	0.48
28:BF:60:LEU:O	28:BF:64:THR:HG22	2.14	0.48
42:BT:30:VAL:HG21	42:BT:79:ALA:HB3	1.94	0.48
47:DY:10:LEU:O	47:DY:13:ALA:HB3	2.13	0.48
41:DS:24:ILE:HG21	41:DS:36:LEU:CD2	2.37	0.48
23:BA:675:A:H4'	27:BE:67:GLN:HE21	1.76	0.48
28:BF:55:LYS:O	28:BF:59:GLU:HG3	2.14	0.48
7:CG:106:GLN:O	7:CG:110:GLN:HG3	2.13	0.48
2:AB:187:LEU:HD11	2:AB:204:ASN:O	2.12	0.48
1:CA:1329:A:P	13:CM:28:ALA:HB3	2.53	0.48
3:AC:30:ARG:CD	14:AN:38:GLY:HA3	2.40	0.48
11:CK:21:ILE:HB	11:CK:84:VAL:HG12	1.95	0.48
37:BO:89:ARG:O	37:BO:90:GLY:O	2.31	0.48
23:DA:2741:A:H2'	23:DA:2742:C:O4'	2.13	0.48
39:DQ:62:ILE:HD12	39:DQ:76:TYR:CE1	2.49	0.48
23:DA:1614:A:C6	41:DS:87:PRO:HA	2.48	0.48
23:BA:478:A:C6	23:BA:480:A:C6	3.01	0.48
23:BA:1478:G:H2'	23:BA:1479:G:H8	1.78	0.48
23:DA:363(A):G:H2'	23:DA:363(B):A:H8	1.78	0.48
23:DA:2815:C:O2'	50:D2:43:HIS:CD2	2.64	0.48
23:DA:826:U:C5	23:DA:828:U:H1'	2.49	0.48
3:AC:91:LEU:HB3	3:AC:99:VAL:HG11	1.94	0.48
2:CB:118:LEU:O	2:CB:122:PHE:HB2	2.13	0.48
23:DA:1995:U:H3'	23:DA:1996:C:H2'	1.96	0.48
1:AA:363:A:H8	12:AL:32:ARG:HH21	1.59	0.48
17:AQ:21:VAL:HG11	17:AQ:59:ILE:HD11	1.95	0.48
1:AA:1154:G:H2'	1:AA:1155:G:C8	2.48	0.48
1:CA:1305:G:H1'	1:CA:1306:A:N7	2.28	0.48
1:AA:277:C:H5''	17:AQ:68:ARG:NH2	2.28	0.48
36:DN:103:ARG:HH12	36:DN:110:PRO:HG3	1.79	0.48
25:BC:259:THR:O	25:BC:260:ARG:HB2	2.14	0.48
23:BA:2887:U:H2'	23:BA:2888:C:H6	1.78	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DA:2590:A:C2	23:DA:2605:U:C2	3.02	0.48
39:BQ:60:LEU:HD23	39:BQ:60:LEU:C	2.33	0.48
1:AA:1366:C:H2'	1:AA:1367:C:C6	2.48	0.48
34:DL:36:LYS:HG3	34:DL:41:ARG:CB	2.40	0.48
23:DA:956:G:N2	23:DA:959:A:H3'	2.29	0.48
1:AA:941:G:C2	1:AA:942:G:C8	3.01	0.48
11:AK:44:SER:OG	11:AK:47:VAL:HG23	2.14	0.48
28:BF:109:VAL:C	28:BF:112:PRO:HD2	2.34	0.48
48:DZ:8:LEU:HA	48:DZ:54:VAL:HG12	1.96	0.48
23:BA:1509:A:H4'	23:BA:1510:A:C1'	2.44	0.48
30:BH:88:ILE:CG2	30:BH:89:TYR:N	2.76	0.48
33:DK:71:ARG:HH12	38:DP:74:ARG:NH2	2.08	0.48
23:DA:2286:A:H4'	23:DA:2287:A:O4'	2.14	0.48
28:BF:98:ARG:O	28:BF:101:ILE:HG12	2.13	0.48
48:BZ:40:THR:O	48:BZ:44:ARG:HG3	2.13	0.48
23:DA:481:G:H1'	23:DA:506:G:N2	2.27	0.48
23:DA:481:G:O2'	23:DA:482:A:P	2.71	0.48
18:AR:56:THR:HB	18:AR:58:LEU:HD13	1.95	0.48
47:BY:52:ASP:O	47:BY:56:GLN:HB2	2.14	0.48
23:DA:768:G:O2'	23:DA:1379:A:N6	2.46	0.48
41:BS:12:ILE:HD12	41:BS:46:PHE:CE2	2.49	0.48
1:AA:269:C:H2'	1:AA:270:A:H8	1.78	0.48
8:AH:11:THR:HG22	8:AH:15:ASN:ND2	2.28	0.48
23:BA:1797:C:O2'	25:BC:259:THR:HG23	2.14	0.48
32:BJ:26:THR:HG22	32:BJ:27:TYR:N	2.29	0.48
23:DA:1051:G:H2'	23:DA:1052:C:C6	2.49	0.48
1:AA:924:C:H2'	1:AA:925:G:H8	1.78	0.48
44:BV:10:ARG:HB3	44:BV:36:LYS:HB3	1.95	0.48
33:DK:12:ASP:OD1	33:DK:85:VAL:HG13	2.12	0.48
13:AM:45:VAL:O	13:AM:48:LEU:HD22	2.13	0.48
23:BA:2335:A:H2'	37:BO:13:ARG:HH22	1.78	0.48
23:DA:1289:C:H2'	23:DA:1290:C:H6	1.79	0.48
1:CA:112:G:C2	1:CA:113:G:C8	3.02	0.48
13:CM:45:VAL:O	13:CM:48:LEU:HD22	2.14	0.48
33:BK:88:ASN:OD1	33:BK:89:ASN:N	2.47	0.48
23:BA:1615:C:O2'	23:BA:1616:A:H5'	2.13	0.48
11:AK:92:GLU:O	11:AK:96:ARG:HG2	2.13	0.48
23:DA:747:U:O2	23:DA:2014:A:H1'	2.13	0.48
34:DL:85:LEU:HA	34:DL:88:LEU:CB	2.43	0.48
43:BU:2:ARG:C	43:BU:4:LYS:H	2.17	0.48
38:DP:46:GLU:OE2	38:DP:89:VAL:HG11	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DA:948:G:OP1	23:DA:962:G:OP1	2.32	0.48
53:B5:53:PRO:HB2	53:B5:57:ARG:NH2	2.28	0.48
1:CA:321:A:C2	1:CA:333:G:C2	3.01	0.48
23:DA:2517:C:C5	23:DA:2542:A:C2	3.02	0.48
2:AB:28:PHE:CD2	2:AB:194:PRO:HG3	2.49	0.48
39:BQ:62:ILE:HD12	39:BQ:76:TYR:CE1	2.48	0.48
4:AD:121:VAL:O	4:AD:134:ASP:HA	2.14	0.48
30:BH:86:THR:O	30:BH:122:GLU:HG3	2.14	0.48
1:CA:404:U:H2'	1:CA:405:U:C6	2.49	0.48
22:CV:6157:A:N6	22:CV:6172:U:H3	2.10	0.48
23:BA:1478:G:O2'	23:BA:1558:A:C2	2.66	0.48
20:AT:26:ASN:HD22	20:AT:27:LYS:H	1.61	0.48
27:BE:28:ILE:O	27:BE:30:PRO:HD3	2.13	0.48
28:BF:128:ARG:HH21	28:BF:130:ASN:HD21	1.61	0.48
1:CA:57:G:H2'	1:CA:58:C:H6	1.77	0.48
38:DP:58:ASN:HD22	38:DP:58:ASN:C	2.17	0.48
27:DE:126:VAL:O	27:DE:196:LEU:HG	2.13	0.48
23:DA:920:G:H2'	23:DA:921:G:H8	1.79	0.48
23:DA:2090:G:H21	46:DX:45:ASN:ND2	2.10	0.48
44:DV:10:ARG:HB3	44:DV:36:LYS:HB3	1.94	0.48
44:BV:10:ARG:HH21	44:BV:26:GLY:H	1.62	0.48
23:DA:1040:C:H2'	23:DA:1041:C:C6	2.49	0.48
24:DB:46:A:H2'	24:DB:47:C:C6	2.48	0.48
23:DA:1239:G:O2'	23:DA:1240:U:H5'	2.13	0.48
23:DA:2366:A:H2'	23:DA:2367:G:O4'	2.14	0.48
32:DJ:160:LYS:CD	32:DJ:161:LEU:H	2.27	0.48
23:DA:754:C:H2'	23:DA:755:C:C6	2.48	0.48
35:DM:23:GLY:HA3	35:DM:98:LYS:HB2	1.95	0.48
29:DG:30:LYS:HB2	29:DG:79:VAL:HA	1.94	0.48
1:AA:562:C:N4	1:AA:884:U:C6	2.82	0.48
23:BA:1399:C:O2'	23:BA:1400:G:H5'	2.14	0.48
28:DF:143:GLU:CD	28:DF:143:GLU:H	2.17	0.48
23:BA:1887:C:H3'	23:BA:1888:G:H5''	1.96	0.48
23:BA:792:G:H5''	23:BA:793:A:H5'	1.94	0.48
23:DA:1689:A:H62	23:DA:1698:A:H2	1.60	0.48
40:DR:49:THR:HB	40:DR:50:PRO:CD	2.43	0.48
1:CA:523:A:N1	12:CL:91:ASP:HB2	2.28	0.48
28:DF:60:LEU:O	28:DF:64:THR:HG22	2.12	0.48
43:DU:81:LYS:CD	43:DU:97:ARG:HB3	2.39	0.48
25:DC:25:THR:HG21	25:DC:81:ALA:CA	2.43	0.48
1:AA:949:A:OP1	13:AM:101:GLN:HB3	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:BC:25:THR:HG21	25:BC:81:ALA:HA	1.95	0.48
23:BA:2027:G:H2'	23:BA:2028:U:O4'	2.14	0.48
12:AL:30:PRO:HB2	12:AL:31:PHE:CD1	2.49	0.48
1:AA:199:G:C2'	1:AA:200:G:H5''	2.41	0.48
3:CC:23:TYR:CG	3:CC:24:ALA:N	2.82	0.48
2:AB:88:ALA:HA	2:AB:223:ILE:HD11	1.95	0.48
36:BN:17:ARG:O	36:BN:20:LEU:HB3	2.14	0.48
19:CS:34:TRP:CZ2	19:CS:57:HIS:HE1	2.32	0.48
1:AA:716:A:N3	11:AK:118:GLY:HA2	2.29	0.48
23:DA:2484:G:H5''	35:DM:45:GLN:HB2	1.95	0.48
23:BA:661:C:H4'	34:BL:16:ARG:HD2	1.95	0.48
41:DS:14:PRO:O	41:DS:15:ARG:C	2.52	0.48
5:AE:81:GLU:HG2	5:AE:90:VAL:HG22	1.96	0.48
8:CH:64:LYS:HD2	8:CH:79:VAL:HG11	1.96	0.48
23:DA:1478:G:N3	23:DA:1479:G:C8	2.82	0.48
46:DX:19:GLN:HA	46:DX:41:ARG:HA	1.96	0.48
23:DA:1390:U:O2'	23:DA:1391:U:H5'	2.13	0.48
23:BA:2061:G:H5''	23:BA:2503:A:C2	2.48	0.48
11:AK:124:LYS:C	11:AK:126:ARG:H	2.16	0.48
32:DJ:80:ALA:HB3	32:DJ:147:ALA:HB2	1.94	0.48
23:DA:2892:A:N6	23:DA:2893:G:C2	2.81	0.48
23:DA:1357:U:H2'	23:DA:1358:G:O4'	2.14	0.48
23:BA:2815:C:O2'	50:B2:43:HIS:CD2	2.66	0.48
1:CA:269:C:H2'	1:CA:270:A:H8	1.78	0.48
1:CA:1281:U:HO2'	1:CA:1282:C:P	2.36	0.48
23:BA:2572:A:H62	26:BD:145:LYS:HG3	1.79	0.48
1:AA:1281:U:HO2'	1:AA:1282:C:P	2.37	0.48
5:CE:11:ILE:HG21	5:CE:105:VAL:HG22	1.96	0.48
8:AH:36:LEU:HA	8:AH:39:LEU:HB3	1.95	0.48
1:AA:1305:G:H1'	1:AA:1306:A:C8	2.48	0.48
23:DA:380:U:O2'	46:DX:20:ARG:HB3	2.13	0.48
1:AA:444:C:H2'	1:AA:445:G:H8	1.78	0.48
23:BA:903:C:H2'	23:BA:904:C:H6	1.79	0.48
1:CA:444:C:H2'	1:CA:445:G:H8	1.79	0.48
23:DA:2784:C:H2'	23:DA:2785:C:H6	1.79	0.48
23:BA:2537:U:H2'	23:BA:2538:C:C6	2.49	0.48
23:BA:1615:C:C6	23:BA:1617:C:C5	3.01	0.48
2:CB:145:LEU:O	2:CB:149:LEU:HB2	2.13	0.48
45:DW:50:ASN:HD22	45:DW:83:PRO:HD3	1.77	0.48
23:BA:649:G:H2'	23:BA:650:C:C6	2.49	0.48
23:DA:1709:U:C2	23:DA:1750:G:C2	3.01	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
37:DO:85:VAL:HG11	37:DO:106:ARG:HD2	1.95	0.48
1:CA:1213:A:O2'	1:CA:1214:C:H5'	2.14	0.48
23:DA:2840:C:H4'	36:DN:53:HIS:CD2	2.49	0.48
53:D5:11:LYS:C	53:D5:11:LYS:HD3	2.34	0.48
41:BS:62:HIS:O	41:BS:64:MET:HG3	2.14	0.48
23:BA:855:G:H2'	23:BA:856:C:C6	2.48	0.48
1:AA:1508:G:H2'	1:AA:1509:C:C6	2.48	0.48
2:CB:52:GLU:O	2:CB:56:ARG:HG3	2.13	0.48
24:BB:78:A:C2	24:BB:99:A:C4	3.02	0.48
44:DV:119:GLU:HG3	44:DV:119:GLU:O	2.14	0.48
23:DA:196:A:H2'	23:DA:196:A:N3	2.28	0.48
1:AA:15:G:C4	1:AA:16:A:C8	3.01	0.48
25:DC:105:ILE:CG1	25:DC:106:ILE:HD12	2.43	0.48
23:DA:2056:G:N2	23:DA:2057:A:C1'	2.77	0.48
42:DT:54:VAL:C	42:DT:55:ASN:HD22	2.17	0.48
12:CL:44:PRO:HD2	12:CL:49:SER:HA	1.94	0.48
53:D5:33:ASN:ND2	53:D5:34:TRP:H	2.11	0.48
1:AA:738:C:H2'	1:AA:739:C:C6	2.48	0.48
35:BM:81:VAL:HG12	35:BM:82:ARG:CG	2.41	0.48
47:BY:10:LEU:O	47:BY:13:ALA:HB3	2.13	0.48
1:AA:675:A:H1'	11:AK:116:HIS:CD2	2.49	0.48
45:BW:32:ARG:CB	45:BW:35:ASN:HD21	2.27	0.48
25:DC:25:THR:O	25:DC:25:THR:HG23	2.14	0.48
48:BZ:8:LEU:HA	48:BZ:54:VAL:HG12	1.95	0.48
25:BC:94:LEU:HB2	25:BC:104:TYR:CE1	2.47	0.48
2:AB:187:LEU:HD23	2:AB:201:ILE:HG22	1.95	0.48
23:BA:1210:A:C5'	23:BA:1210:A:H8	2.22	0.48
2:AB:28:PHE:CD1	2:AB:190:THR:HA	2.48	0.48
23:BA:310:A:OP1	43:BU:17:SER:O	2.31	0.48
28:DF:77:ILE:HG22	28:DF:80:PHE:N	2.25	0.48
48:DZ:40:THR:O	48:DZ:44:ARG:HG3	2.14	0.48
37:DO:39:ILE:O	37:DO:48:LEU:HD13	2.14	0.48
29:BG:46:GLU:HG3	29:BG:51:ARG:NE	2.28	0.48
25:DC:79:VAL:O	25:DC:113:VAL:HG13	2.14	0.48
29:BG:58:GLU:HB2	29:BG:61:HIS:ND1	2.29	0.48
34:DL:95:VAL:HG23	34:DL:125:VAL:HA	1.95	0.48
23:BA:2244:U:O2'	23:BA:2245:U:H5'	2.13	0.48
1:AA:32:A:C6	1:AA:33:A:C6	3.02	0.48
1:CA:820:U:H4'	1:CA:821:G:OP2	2.14	0.48
44:DV:5:LEU:HD23	44:DV:6:LYS:N	2.29	0.48
1:AA:363:A:C8	12:AL:32:ARG:NH2	2.82	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:600:C:OP1	8:AH:97:VAL:HG12	2.14	0.48
23:BA:1639:U:H4'	23:BA:2699:C:H4'	1.94	0.48
1:AA:18:C:H5''	5:AE:127:ASN:ND2	2.28	0.48
23:DA:380:U:H1'	46:DX:20:ARG:NH1	2.29	0.48
4:AD:92:VAL:O	4:AD:96:LEU:HD23	2.14	0.48
23:BA:2334:G:H4'	23:BA:2335:A:OP2	2.14	0.48
1:CA:705:U:C5	1:CA:706:A:C5	3.02	0.48
10:AJ:3:LYS:HD2	10:AJ:77:PRO:HD3	1.96	0.48
45:DW:66:VAL:O	45:DW:81:VAL:HA	2.13	0.48
23:DA:714:U:O2	23:DA:716:A:C8	2.67	0.48
5:CE:70:PRO:HB3	5:CE:144:THR:HG22	1.95	0.48
30:DH:66:GLU:HB3	30:DH:67:ARG:NH1	2.28	0.48
25:DC:40:THR:HG22	25:DC:41:GLY:N	2.28	0.48
5:CE:76:ILE:HG23	5:CE:78:HIS:H	1.78	0.48
43:BU:9:LYS:O	43:BU:27:VAL:HG21	2.14	0.48
4:CD:51:PRO:HB3	4:CD:55:ALA:HB3	1.96	0.48
23:BA:1517:G:H2'	23:BA:1518:C:C6	2.49	0.48
6:AF:69:GLU:CD	6:AF:69:GLU:H	2.16	0.48
5:CE:81:GLU:HG2	5:CE:90:VAL:HG22	1.96	0.48
23:BA:242:G:N7	53:B5:5:LYS:HG2	2.28	0.48
1:AA:624:C:H2'	1:AA:625:G:H8	1.78	0.48
23:BA:2850:A:H2'	23:BA:2851:A:C8	2.49	0.48
1:AA:1327:C:H2'	1:AA:1328:C:H6	1.78	0.48
23:DA:1504:C:O2'	23:DA:1505:C:O5'	2.31	0.48
1:AA:1017:G:H2'	1:AA:1018:C:H6	1.79	0.48
1:CA:688:G:H2'	1:CA:689:C:H6	1.78	0.48
24:BB:8:U:H5''	37:BO:15:ARG:NH2	2.29	0.48
36:DN:87:TYR:HE1	36:DN:117:VAL:HG13	1.79	0.48
25:DC:72:LYS:HE3	25:DC:101:GLU:CB	2.44	0.48
5:CE:31:LEU:HD23	5:CE:32:VAL:N	2.29	0.48
27:BE:37:VAL:HG13	27:BE:184:TYR:HD1	1.78	0.48
23:DA:270(Q):C:O2'	23:DA:270(R):C:H6	1.97	0.48
23:DA:356:G:H2'	23:DA:357:A:C8	2.49	0.48
38:BP:24:PRO:HA	38:BP:49:VAL:HG13	1.95	0.48
24:DB:28:C:H2'	24:DB:29:A:O4'	2.14	0.48
32:BJ:77:VAL:HB	32:BJ:145:VAL:HG22	1.94	0.48
39:DQ:107:ALA:O	39:DQ:111:GLU:HG2	2.13	0.48
4:AD:93:PHE:O	4:AD:97:LEU:HG	2.14	0.48
53:D5:29:LYS:NZ	53:D5:29:LYS:HB3	2.29	0.48
45:DW:27:GLU:HB2	45:DW:69:PHE:HD1	1.79	0.48
31:BI:9:LEU:O	31:BI:13:LEU:HG	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:AE:76:ILE:HG23	5:AE:78:HIS:H	1.78	0.48
23:DA:954:G:C5	23:DA:955:C:C5	3.01	0.48
1:AA:92:G:H2'	1:AA:93:U:O4'	2.14	0.48
1:AA:91:C:O2'	1:AA:92:G:H5'	2.14	0.48
17:CQ:54:GLY:HA3	17:CQ:82:MET:CE	2.44	0.48
36:BN:50:HIS:O	36:BN:54:LEU:HB2	2.14	0.48
38:BP:27:THR:CG2	38:BP:90:GLN:HB3	2.44	0.48
1:CA:675:A:H2'	1:CA:676:A:H8	1.79	0.48
1:CA:941:G:C2	1:CA:942:G:C8	3.02	0.48
6:AF:19:LEU:O	6:AF:23:LYS:HG3	2.13	0.48
25:BC:154:LYS:C	25:BC:155:LEU:HD12	2.34	0.48
37:DO:34:HIS:CG	37:DO:54:LEU:HB2	2.49	0.48
13:AM:3:ARG:NH1	28:BF:113:ARG:HD2	2.28	0.48
30:DH:109:ILE:N	30:DH:109:ILE:HD13	2.29	0.48
1:AA:404:U:H2'	1:AA:405:U:C6	2.49	0.48
8:AH:48:TYR:HA	8:AH:60:ARG:O	2.14	0.48
23:BA:2599:G:N7	25:BC:237:GLU:HG3	2.29	0.48
36:DN:97:VAL:HA	36:DN:113:LEU:O	2.14	0.48
1:CA:1349:A:H2'	1:CA:1350:A:O4'	2.14	0.48
23:BA:39:C:H2'	23:BA:40:C:C6	2.49	0.48
19:AS:5:LEU:HG	19:AS:10:PHE:HB3	1.95	0.48
1:AA:10:A:H2'	1:AA:11:G:H8	1.78	0.48
1:CA:1504:G:O2'	1:CA:1505:G:P	2.72	0.48
40:BR:81:TYR:O	40:BR:82:ARG:HG3	2.13	0.48
38:BP:60:THR:HG22	38:BP:77:PRO:HA	1.96	0.48
1:AA:1070:U:H2'	1:AA:1071:C:C6	2.48	0.48
1:CA:622:A:C8	1:CA:623:C:C6	3.02	0.48
44:DV:81:ARG:O	44:DV:82:ARG:HB2	2.13	0.48
8:CH:36:LEU:HA	8:CH:39:LEU:HB3	1.96	0.48
23:DA:1973:G:H2'	23:DA:1974:C:C6	2.49	0.48
23:DA:2252:G:H2'	23:DA:2253:G:C8	2.49	0.48
23:DA:2537:U:H2'	23:DA:2538:C:C6	2.48	0.48
1:CA:9:G:H2'	1:CA:10:A:H8	1.79	0.48
17:CQ:76:LEU:HD12	17:CQ:77:VAL:H	1.78	0.48
1:AA:1305:G:H1'	1:AA:1306:A:N7	2.28	0.48
8:AH:82:HIS:HD2	8:AH:138:TRP:NE1	2.12	0.48
1:AA:1084:G:H5'	1:AA:1102:A:OP2	2.14	0.48
30:BH:117:GLU:HG3	30:BH:118:LYS:N	2.28	0.48
2:AB:61:LEU:HD21	2:AB:68:ILE:HG12	1.95	0.48
25:BC:40:THR:HG22	25:BC:41:GLY:N	2.28	0.48
4:AD:51:PRO:HB3	4:AD:55:ALA:HB3	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DA:649:G:H2'	23:DA:650:C:C6	2.49	0.48
24:BB:6:C:C2	24:BB:115:G:N2	2.82	0.48
1:CA:857:C:H2'	1:CA:858:G:O4'	2.14	0.48
23:BA:409:C:O2'	23:BA:410:G:H5'	2.13	0.48
50:D2:3:LYS:O	50:D2:4:HIS:C	2.52	0.47
23:DA:587:C:C5	23:DA:671:C:H1'	2.49	0.47
40:DR:47:VAL:HG12	40:DR:49:THR:O	2.13	0.47
15:CO:33:THR:HG23	15:CO:63:ARG:HH22	1.79	0.47
23:BA:114(B):A:C4	23:BA:1144:G:C8	3.01	0.47
23:DA:761:A:O5'	23:DA:761:A:C8	2.48	0.47
42:DT:30:VAL:HG21	42:DT:79:ALA:HB3	1.96	0.47
43:BU:76:CYS:CB	43:BU:77:PRO:CD	2.92	0.47
38:DP:27:THR:HA	38:DP:48:ILE:HA	1.96	0.47
53:B5:14:VAL:CG1	53:B5:22:VAL:HG13	2.43	0.47
23:BA:124:G:N7	52:B4:19:ARG:NH2	2.62	0.47
43:DU:10:GLY:HA2	43:DU:27:VAL:HG23	1.95	0.47
26:DD:101:ARG:HG2	26:DD:171:GLU:HA	1.96	0.47
26:DD:171:GLU:HG2	26:DD:185:LYS:HG2	1.96	0.47
16:CP:18:ARG:HD3	16:CP:35:LYS:HD2	1.95	0.47
30:BH:79:ILE:HG22	30:BH:81:VAL:CG2	2.44	0.47
23:BA:603:A:C6	23:BA:655:A:H1'	2.48	0.47
2:AB:20:GLU:HA	2:AB:20:GLU:OE1	2.13	0.47
10:AJ:34:VAL:HG22	10:AJ:74:ILE:HG22	1.96	0.47
23:BA:2286:A:H4'	23:BA:2287:A:O4'	2.14	0.47
23:DA:1317:A:N6	23:DA:1336:A:N6	2.62	0.47
23:BA:661:C:H4'	34:BL:16:ARG:CD	2.44	0.47
34:BL:16:ARG:CZ	34:BL:18:ARG:HG3	2.44	0.47
37:DO:89:ARG:O	37:DO:90:GLY:O	2.31	0.47
9:AI:14:VAL:O	9:AI:65:VAL:HG23	2.14	0.47
9:CI:62:TYR:C	9:CI:63:ILE:HD12	2.35	0.47
23:BA:1478:G:N3	23:BA:1479:G:C8	2.82	0.47
23:DA:598:G:H5'	34:DL:15:ARG:CB	2.44	0.47
9:CI:99:LEU:HD12	9:CI:101:PHE:HE2	1.78	0.47
1:AA:1349:A:H2'	1:AA:1350:A:H8	1.76	0.47
23:DA:226:G:N2	23:DA:228:A:H62	2.12	0.47
18:CR:36:ASN:HD22	18:CR:39:VAL:HG21	1.78	0.47
5:AE:101:ILE:HD11	5:AE:119:LEU:CD2	2.44	0.47
23:DA:2476:A:C2	23:DA:2477:C:C6	3.02	0.47
23:BA:784:A:N7	25:BC:229:VAL:HG21	2.29	0.47
34:DL:135:LEU:HD13	34:DL:135:LEU:O	2.14	0.47
38:DP:60:THR:HG22	38:DP:77:PRO:HA	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
44:BV:5:LEU:HD23	44:BV:6:LYS:N	2.29	0.47
1:AA:963:G:H2'	1:AA:964:A:H8	1.77	0.47
1:CA:865:A:O5'	1:CA:865:A:H8	1.97	0.47
10:AJ:80:LYS:HB2	10:AJ:80:LYS:NZ	2.30	0.47
30:BH:25:TYR:O	30:BH:29:TYR:HB3	2.14	0.47
1:AA:1203:C:H2'	1:AA:1204:A:C8	2.49	0.47
41:BS:84:ARG:HB2	41:BS:96:ILE:HG22	1.96	0.47
23:BA:1796:U:H2'	23:BA:1797:C:C6	2.49	0.47
1:AA:832:C:N4	1:AA:855:G:O6	2.47	0.47
1:CA:1455:G:O2'	20:CT:28:ALA:HB1	2.14	0.47
18:CR:43:PHE:O	18:CR:51:LEU:HB2	2.13	0.47
8:CH:23:SER:HB3	8:CH:62:TYR:HA	1.96	0.47
23:DA:909:A:H2'	23:DA:912:C:H5	1.79	0.47
23:DA:1198:U:O2	23:DA:1249:U:H1'	2.14	0.47
26:BD:31:CYS:HB3	26:BD:49:LEU:HB3	1.95	0.47
32:DJ:63:PRO:O	39:DQ:64:ARG:HD2	2.14	0.47
23:DA:409:C:O2'	23:DA:410:G:H5'	2.14	0.47
40:BR:22:VAL:CG1	40:BR:23:GLU:N	2.76	0.47
23:DA:1971:A:C5	25:DC:241:PRO:HG3	2.48	0.47
2:CB:162:ILE:HD11	2:CB:184:VAL:HG13	1.96	0.47
28:BF:92:VAL:O	28:BF:92:VAL:HG13	2.14	0.47
28:DF:92:VAL:HG13	28:DF:92:VAL:O	2.15	0.47
23:BA:827:U:O2	23:BA:2246:G:H4'	2.14	0.47
23:BA:783:A:C3'	23:BA:783:A:C8	2.97	0.47
35:BM:81:VAL:HG12	35:BM:82:ARG:N	2.29	0.47
16:AP:18:ARG:HD3	16:AP:35:LYS:HD2	1.96	0.47
26:DD:5:LEU:HB2	26:DD:51:PHE:HD2	1.77	0.47
23:DA:603:A:C6	23:DA:655:A:H1'	2.50	0.47
19:CS:28:LYS:HB3	19:CS:29:ARG:NH1	2.29	0.47
17:AQ:66:SER:OG	17:AQ:69:LYS:HB3	2.13	0.47
23:BA:1448:G:H21	23:BA:1529:A:H2	1.62	0.47
23:BA:861:A:H2'	23:BA:862:G:O4'	2.15	0.47
40:DR:79:VAL:O	40:DR:79:VAL:HG12	2.14	0.47
3:AC:18:TRP:HE3	3:AC:18:TRP:H	1.61	0.47
23:BA:1331:A:O2'	23:BA:1332:G:C8	2.66	0.47
20:CT:26:ASN:HB2	20:CT:71:THR:CG2	2.44	0.47
30:DH:6:LEU:HD23	30:DH:36:ALA:HA	1.96	0.47
23:DA:1308:A:N6	23:DA:1309:G:C2	2.83	0.47
23:DA:2593:U:H2'	23:DA:2594:C:H6	1.79	0.47
5:CE:101:ILE:HD11	5:CE:119:LEU:CD2	2.44	0.47
1:AA:1062:U:H2'	1:AA:1063:C:C6	2.49	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
30:DH:57:ARG:O	30:DH:61:ARG:HG3	2.14	0.47
4:CD:79:PHE:CE1	4:CD:204:ILE:HA	2.49	0.47
1:AA:820:U:H4'	1:AA:821:G:OP2	2.13	0.47
8:CH:82:HIS:HD2	8:CH:138:TRP:NE1	2.12	0.47
16:CP:45:THR:HB	16:CP:46:PRO:HD2	1.96	0.47
23:BA:634:C:H2'	23:BA:635:C:C6	2.50	0.47
23:DA:2552:U:H2'	23:DA:2554:U:OP2	2.14	0.47
20:CT:14:LYS:O	20:CT:18:GLN:HG3	2.13	0.47
23:BA:471:A:H2'	23:BA:472:A:O4'	2.14	0.47
23:BA:1289:C:H2'	23:BA:1290:C:H6	1.79	0.47
41:BS:9:TYR:H	41:BS:102:HIS:HD2	1.63	0.47
1:CA:746:A:H2'	1:CA:747:C:C6	2.50	0.47
23:BA:2512:C:H5''	23:BA:2513:G:OP2	2.14	0.47
23:BA:1040:C:H2'	23:BA:1041:C:C6	2.49	0.47
45:DW:12:ASN:O	45:DW:14:ARG:HB2	2.14	0.47
23:BA:2751:G:H2'	23:BA:2751:G:N3	2.29	0.47
23:BA:239:U:O2'	23:BA:240:G:H5'	2.14	0.47
1:AA:171:A:H2'	1:AA:172:A:C8	2.49	0.47
1:CA:1508:G:H2'	1:CA:1509:C:C6	2.49	0.47
29:DG:154:PRO:HB3	29:DG:163:TYR:CE2	2.49	0.47
1:AA:636:U:H5'	17:AQ:2:PRO:HD3	1.95	0.47
12:CL:44:PRO:CD	12:CL:50:ALA:H	2.28	0.47
12:CL:26:LEU:HD13	12:CL:27:LYS:N	2.18	0.47
21:CU:22:ARG:HD2	21:CU:23:PRO:CD	2.37	0.47
1:CA:738:C:H2'	1:CA:739:C:C6	2.48	0.47
25:BC:136:ILE:O	25:BC:168:ARG:NH2	2.47	0.47
10:AJ:33:GLN:HB2	10:AJ:75:ILE:HD11	1.96	0.47
23:DA:2822:G:O5'	23:DA:2822:G:H8	1.97	0.47
23:BA:124:G:C5	52:B4:19:ARG:NH2	2.82	0.47
12:AL:76:LEU:HD11	12:AL:106:ALA:HA	1.96	0.47
26:DD:103:ASP:OD1	26:DD:201:THR:HG23	2.15	0.47
16:CP:22:THR:HG22	16:CP:32:TYR:HA	1.95	0.47
23:DA:1509:A:H4'	23:DA:1510:A:C1'	2.44	0.47
23:DA:2542:A:C8	23:DA:2544:G:O6	2.67	0.47
10:CJ:34:VAL:CG2	10:CJ:74:ILE:HG22	2.43	0.47
30:DH:78:THR:HA	30:DH:143:SER:CB	2.41	0.47
23:BA:1504:C:O2'	23:BA:1505:C:O5'	2.31	0.47
5:CE:91:LEU:HD12	5:CE:91:LEU:N	2.30	0.47
12:CL:24:PRO:HD2	12:CL:97:TYR:OH	2.14	0.47
1:AA:370:C:H2'	1:AA:371:G:C8	2.47	0.47
1:AA:114:U:H2'	1:AA:115:G:H8	1.75	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BA:966:G:C6	23:BA:967:C:N4	2.82	0.47
23:BA:768:G:O2'	23:BA:1379:A:N6	2.47	0.47
1:AA:255:G:H2'	1:AA:256:U:H6	1.78	0.47
23:DA:919:G:H2'	23:DA:920:G:C8	2.50	0.47
36:BN:96:ARG:NH2	36:BN:117:VAL:HG23	2.29	0.47
5:AE:31:LEU:HD23	5:AE:32:VAL:N	2.29	0.47
35:BM:47:ILE:HG22	35:BM:48:GLU:N	2.28	0.47
1:CA:1423:G:H2'	1:CA:1424:C:H6	1.79	0.47
23:DA:903:C:H2'	23:DA:904:C:H6	1.79	0.47
26:DD:78:LEU:N	26:DD:78:LEU:HD23	2.29	0.47
1:CA:858:G:O6	1:CA:869:G:C8	2.67	0.47
23:BA:714:U:O2	23:BA:716:A:C8	2.67	0.47
23:DA:1411:C:H2'	23:DA:1412:A:C8	2.50	0.47
1:AA:1480:G:C5	1:AA:1481:U:C5	3.03	0.47
1:AA:1213:A:O2'	1:AA:1214:C:H5'	2.14	0.47
23:BA:1871:A:H2'	23:BA:1872:A:C8	2.50	0.47
40:DR:99:ILE:N	40:DR:99:ILE:HD13	2.28	0.47
23:DA:2056:G:C2'	23:DA:2056:G:N3	2.77	0.47
1:AA:1240:U:OP1	7:AG:115:ARG:HA	2.15	0.47
35:DM:81:VAL:HG12	35:DM:82:ARG:N	2.29	0.47
37:BO:69:VAL:O	37:BO:72:ALA:HB3	2.14	0.47
36:BN:54:LEU:HD23	36:BN:62:ALA:HB1	1.97	0.47
25:DC:25:THR:O	25:DC:27:THR:N	2.48	0.47
26:BD:171:GLU:HG2	26:BD:185:LYS:HG2	1.96	0.47
23:DA:2038:G:H2'	23:DA:2039:C:H6	1.78	0.47
23:BA:2516:G:C6	23:BA:2517:C:N4	2.83	0.47
3:AC:11:ARG:HH21	3:AC:180:ALA:HB3	1.80	0.47
2:AB:47:THR:O	2:AB:51:LEU:HG	2.14	0.47
26:BD:51:PHE:HB3	26:BD:52:LEU:HD12	1.96	0.47
25:DC:69:ARG:HH12	25:DC:117:VAL:CG2	2.26	0.47
23:DA:94:G:H21	47:DY:47:ASN:ND2	2.12	0.47
23:DA:1478:G:O2'	23:DA:1558:A:C2	2.67	0.47
46:DX:19:GLN:HG2	46:DX:41:ARG:CB	2.45	0.47
23:BA:363(A):G:H2'	23:BA:363(B):A:H8	1.79	0.47
35:DM:39:PRO:O	35:DM:40:ALA:HB2	2.14	0.47
35:BM:6:ARG:N	35:BM:6:ARG:HE	2.12	0.47
1:AA:1097:C:H2'	1:AA:1098:C:H6	1.79	0.47
27:BE:126:VAL:O	27:BE:196:LEU:HG	2.14	0.47
30:DH:58:LEU:HD23	30:DH:61:ARG:HD2	1.96	0.47
1:AA:186(A):C:C5'	20:AT:78:ALA:HB1	2.43	0.47
27:DE:125:LEU:HB3	27:DE:196:LEU:HD23	1.95	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DA:2699:C:H2'	23:DA:2700:C:O4'	2.14	0.47
1:CA:1017:G:H2'	1:CA:1018:C:H6	1.79	0.47
23:BA:380:U:H1'	46:BX:20:ARG:NH1	2.29	0.47
36:BN:87:TYR:HE1	36:BN:117:VAL:HG13	1.79	0.47
44:BV:27:VAL:HG22	44:BV:36:LYS:HA	1.95	0.47
23:DA:1709:U:H2'	23:DA:1710:C:C6	2.49	0.47
36:DN:53:HIS:O	36:DN:56:LYS:HB3	2.14	0.47
23:BA:356:G:H2'	23:BA:357:A:C8	2.49	0.47
23:DA:451:C:H41	23:DA:453:C:H3'	1.78	0.47
23:DA:1936:A:H3'	23:DA:1936:A:OP1	2.14	0.47
26:DD:152:LYS:HE2	26:DD:152:LYS:HB3	1.72	0.47
37:BO:85:VAL:HG11	37:BO:106:ARG:HD2	1.95	0.47
50:D2:4:HIS:HB2	50:D2:5:PRO:CD	2.45	0.47
23:DA:1827:C:H2'	23:DA:1828:G:O4'	2.14	0.47
1:CA:522:C:N4	1:CA:528:C:H42	2.11	0.47
10:AJ:49:VAL:HG21	14:AN:41:ARG:CB	2.44	0.47
34:BL:84:ASN:HA	34:BL:115:LEU:O	2.15	0.47
23:DA:124:G:C5	52:D4:19:ARG:NH2	2.82	0.47
23:DA:1578:U:H2'	23:DA:1579:A:H5''	1.95	0.47
7:AG:69:VAL:HG12	7:AG:69:VAL:O	2.15	0.47
3:AC:23:TYR:HA	10:AJ:11:PHE:CE1	2.50	0.47
23:BA:2481:G:HO2'	23:BA:2482:G:P	2.37	0.47
23:DA:1431:U:H2'	23:DA:1432:C:H6	1.80	0.47
30:DH:12:LEU:N	30:DH:12:LEU:HD22	2.29	0.47
19:AS:28:LYS:HB3	19:AS:29:ARG:NH1	2.29	0.47
4:AD:122:ARG:HD3	4:AD:122:ARG:C	2.34	0.47
24:BB:75:G:N1	24:BB:102:G:N2	2.61	0.47
23:DA:2687:U:C4	23:DA:2688:U:H5	2.30	0.47
29:BG:103:LEU:HD22	29:BG:123:PHE:CE1	2.50	0.47
26:BD:4:ILE:HG12	26:BD:28:ALA:HB1	1.95	0.47
23:DA:2630:G:H1'	23:DA:2894:G:H1'	1.97	0.47
11:AK:59:TYR:CZ	11:AK:63:LEU:HD11	2.49	0.47
1:CA:716:A:N3	11:CK:118:GLY:HA2	2.29	0.47
23:DA:320:A:H5''	23:DA:321:G:OP1	2.15	0.47
8:AH:97:VAL:HG13	8:AH:98:LYS:H	1.80	0.47
4:AD:79:PHE:CE1	4:AD:204:ILE:HA	2.49	0.47
13:CM:15:VAL:HG13	13:CM:43:THR:O	2.14	0.47
1:CA:1410:G:H2'	1:CA:1411:C:C6	2.49	0.47
23:DA:2846:G:H2'	23:DA:2847:U:C6	2.50	0.47
28:BF:16:ARG:HB3	28:BF:17:PRO:HD3	1.97	0.47
5:CE:65:ASN:O	5:CE:66:MET:HB2	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:46:G:O2'	1:AA:365:U:H1'	2.15	0.47
23:BA:1341:U:O4	42:BT:16:LYS:HE2	2.14	0.47
25:BC:76:PRO:HB3	25:BC:116:GLN:HE21	1.80	0.47
4:CD:93:PHE:O	4:CD:97:LEU:HG	2.13	0.47
50:B2:18:ALA:O	50:B2:21:SER:HB2	2.14	0.47
23:BA:1709:U:H2'	23:BA:1710:C:C6	2.49	0.47
12:CL:78:GLU:O	12:CL:78:GLU:HG2	2.14	0.47
28:BF:143:GLU:CD	28:BF:143:GLU:H	2.17	0.47
40:BR:99:ILE:N	40:BR:99:ILE:HD13	2.29	0.47
28:BF:104:GLU:O	28:BF:108:ASN:HB2	2.14	0.47
1:AA:1446:A:N1	38:BP:118:ARG:CZ	2.78	0.47
23:BA:1022:G:N2	23:BA:114(B):A:C2	2.83	0.47
23:DA:2247:A:H2'	23:DA:2248:C:C6	2.49	0.47
37:BO:49:VAL:CG1	37:BO:76:LYS:HB2	2.45	0.47
23:DA:1658:C:OP1	26:DD:132:HIS:O	2.33	0.47
3:AC:23:TYR:CG	3:AC:24:ALA:N	2.82	0.47
26:BD:5:LEU:C	26:BD:51:PHE:HE2	2.18	0.47
23:DA:775:G:C5	23:DA:794:G:C8	3.02	0.47
46:BX:27:GLU:HB2	46:BX:32:LYS:O	2.14	0.47
24:DB:106:G:C6	24:DB:107:U:C4	3.03	0.47
1:AA:1194:U:H2'	1:AA:1195:C:H6	1.76	0.47
23:DA:1030:G:OP2	35:DM:128:LYS:HE3	2.14	0.47
26:DD:4:ILE:HG12	26:DD:28:ALA:HB1	1.95	0.47
19:CS:40:ILE:HG21	19:CS:62:ILE:HD11	1.97	0.47
25:DC:108:PRO:CG	25:DC:143:HIS:CE1	2.97	0.47
47:BY:41:ILE:HD11	47:BY:44:LEU:HD12	1.95	0.47
41:DS:4:LYS:HG2	41:DS:106:ILE:CG2	2.44	0.47
24:DB:111:U:HO2'	24:DB:112:G:H8	1.59	0.47
44:BV:81:ARG:O	44:BV:82:ARG:HB2	2.14	0.47
36:DN:87:TYR:OH	36:DN:116:LEU:HB3	2.15	0.47
36:DN:96:ARG:NH2	36:DN:117:VAL:HG23	2.30	0.47
4:CD:163:GLU:O	4:CD:166:LYS:HG3	2.14	0.47
38:DP:109:GLU:HA	38:DP:112:ARG:HG3	1.97	0.47
1:CA:1203:C:H2'	1:CA:1204:A:C8	2.50	0.47
21:CU:6:ARG:HG3	21:CU:15:ARG:NH1	2.30	0.47
23:BA:2836:U:C4	23:BA:2883:A:N6	2.82	0.47
26:DD:54:GLN:HB2	26:DD:74:PRO:O	2.15	0.47
1:CA:617:G:H4'	16:CP:44:THR:HB	1.95	0.47
13:CM:81:LEU:HD22	13:CM:86:CYS:SG	2.54	0.47
1:CA:632:A:H2'	1:CA:633:G:O4'	2.15	0.47
6:CF:82:ARG:HA	6:CF:82:ARG:HH11	1.78	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:375:U:H4'	16:AP:17:TYR:CE2	2.49	0.47
4:CD:195:ALA:C	4:CD:196:LEU:HD12	2.35	0.47
29:BG:154:PRO:HB3	29:BG:163:TYR:CE2	2.49	0.47
14:AN:40:CYS:SG	14:AN:42:ILE:HB	2.55	0.47
23:BA:1326:U:O2'	23:BA:1327:C:H5'	2.15	0.47
23:BA:718:A:O5'	23:BA:718:A:H8	1.96	0.47
23:DA:2887:U:H2'	23:DA:2888:C:H6	1.80	0.47
53:B5:60:LEU:C	53:B5:62:LEU:H	2.17	0.47
39:BQ:92:ARG:HG2	40:BR:11:GLN:CD	2.35	0.47
25:BC:238:GLY:O	25:BC:240:ALA:N	2.48	0.47
15:AO:33:THR:HG23	15:AO:63:ARG:HH22	1.80	0.47
23:DA:1900:A:N1	23:DA:1970:A:C6	2.83	0.47
43:DU:13:VAL:HG13	43:DU:73:ARG:O	2.14	0.47
53:D5:32:LEU:HD23	53:D5:33:ASN:H	1.80	0.47
41:BS:24:ILE:HG21	41:BS:36:LEU:CD2	2.36	0.47
5:AE:51:VAL:HB	5:AE:52:PRO:CD	2.37	0.47
1:CA:91:C:O2'	1:CA:92:G:H5'	2.15	0.47
23:BA:270(L):C:H6	23:BA:270(L):C:O5'	1.97	0.47
38:DP:27:THR:CG2	38:DP:90:GLN:HB3	2.44	0.47
44:BV:99:TYR:CE2	44:BV:125:LEU:HD12	2.49	0.47
36:DN:54:LEU:HD23	36:DN:62:ALA:HB1	1.95	0.47
34:DL:6:LEU:N	34:DL:6:LEU:HD23	2.23	0.47
16:AP:22:THR:HG22	16:AP:32:TYR:HA	1.96	0.47
23:BA:744:G:OP1	26:BD:132:HIS:CB	2.62	0.47
30:BH:83:ALA:CB	30:BH:123:LEU:HD12	2.43	0.47
30:BH:82:ARG:HB3	30:BH:89:TYR:CG	2.49	0.47
2:AB:91:PRO:CA	2:AB:154:LEU:HD11	2.42	0.47
3:CC:23:TYR:HA	10:CJ:11:PHE:CE1	2.50	0.47
23:DA:2469:A:H5'	23:DA:2470:G:OP2	2.15	0.47
30:BH:130:TYR:HD2	30:BH:132:PRO:HG3	1.80	0.47
10:CJ:34:VAL:HG22	10:CJ:74:ILE:HG22	1.96	0.47
1:AA:406:G:N2	1:AA:437:U:C2	2.83	0.47
29:BG:51:ARG:O	29:BG:52:VAL:HG23	2.15	0.47
41:BS:14:PRO:O	41:BS:15:ARG:C	2.52	0.47
3:CC:18:TRP:H	3:CC:18:TRP:HE3	1.62	0.47
1:CA:1073:U:H2'	1:CA:1074:G:C8	2.47	0.47
1:AA:920:U:H2'	1:AA:921:U:H6	1.80	0.47
1:AA:1151:A:O2'	1:AA:1152:A:C8	2.65	0.47
23:BA:2365:G:O6	53:B5:39:LYS:HE3	2.15	0.47
35:DM:6:ARG:N	35:DM:6:ARG:HE	2.13	0.47
32:DJ:127:LYS:HB2	32:DJ:140:PHE:HE1	1.77	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BA:1357:U:H2'	23:BA:1358:G:O4'	2.15	0.47
34:BL:126:VAL:HA	34:BL:145:PRO:HB2	1.95	0.47
1:AA:57:G:H2'	1:AA:58:C:H6	1.77	0.47
23:DA:2564:A:OP1	23:DA:2648:C:H4'	2.14	0.47
46:DX:58:ILE:HD11	46:DX:91:LYS:CG	2.44	0.47
1:AA:1504:G:O2'	1:AA:1505:G:P	2.73	0.47
1:CA:15:G:C4	1:CA:16:A:C8	3.03	0.47
23:DA:1639:U:H4'	23:DA:2699:C:H4'	1.96	0.47
1:AA:818:G:H3'	1:AA:819:A:H5''	1.97	0.47
23:DA:1833:U:C2	23:DA:1834:U:C6	3.01	0.47
1:AA:973:G:H3'	1:AA:974:A:H5''	1.97	0.47
23:BA:1495:A:H2'	23:BA:1496:A:C2	2.50	0.47
23:BA:2730:C:O2'	23:BA:2731:G:H5'	2.14	0.47
23:BA:628:G:H2'	23:BA:629:G:H8	1.79	0.47
2:AB:17:PHE:HB2	2:AB:42:ILE:HG22	1.95	0.47
1:CA:1410:G:H2'	1:CA:1411:C:H6	1.80	0.47
1:AA:1412:C:H2'	1:AA:1413:A:H8	1.79	0.47
23:BA:2784:C:H2'	23:BA:2785:C:H6	1.80	0.47
5:AE:144:THR:O	5:AE:148:VAL:HG23	2.15	0.47
26:BD:49:LEU:O	26:BD:78:LEU:HA	2.15	0.47
23:BA:2836:U:H2'	23:BA:2837:G:C8	2.50	0.47
1:AA:746:A:H2'	1:AA:747:C:C6	2.50	0.47
1:AA:102:G:H2'	1:AA:103:C:C6	2.49	0.47
28:DF:104:GLU:O	28:DF:108:ASN:HB2	2.15	0.47
41:DS:22:ASP:HA	41:DS:25:ARG:HH12	1.79	0.47
23:DA:816:C:O2'	23:DA:817:C:H5'	2.15	0.47
23:DA:1449:G:H2'	23:DA:1450:C:C6	2.50	0.47
3:AC:113:ALA:HB3	3:AC:114:PRO:HD3	1.97	0.47
6:AF:26:ILE:O	6:AF:30:LEU:HG	2.15	0.47
44:BV:119:GLU:HG3	44:BV:119:GLU:O	2.14	0.47
23:BA:1051:G:H2'	23:BA:1052:C:C6	2.49	0.47
23:DA:2751:G:H2'	23:DA:2751:G:N3	2.30	0.47
23:BA:1121:C:O5'	23:BA:1121:C:H6	1.98	0.47
41:DS:62:HIS:O	41:DS:64:MET:HG3	2.14	0.47
11:CK:92:GLU:O	11:CK:96:ARG:HG2	2.13	0.47
45:BW:66:VAL:O	45:BW:81:VAL:HA	2.15	0.47
25:DC:259:THR:O	25:DC:260:ARG:HB2	2.15	0.47
32:BJ:30:LYS:O	32:BJ:32:VAL:HG23	2.15	0.47
7:CG:101:LEU:O	7:CG:105:VAL:HG23	2.15	0.47
41:DS:20:VAL:O	41:DS:23:LEU:HB2	2.14	0.47
32:DJ:30:LYS:O	32:DJ:32:VAL:HG23	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:DD:176:ILE:O	26:DD:176:ILE:HG22	2.15	0.47
1:AA:45:U:H6	1:AA:45:U:O5'	1.97	0.47
1:AA:342:C:N3	1:AA:348:G:C2	2.83	0.47
24:DB:78:A:C2	24:DB:99:A:C4	3.02	0.47
23:BA:828:U:O2	23:BA:828:U:H3'	2.15	0.47
34:DL:50:ARG:HD3	34:DL:51:PHE:HB2	1.97	0.47
40:BR:6:LYS:HA	40:BR:11:GLN:HB3	1.97	0.47
23:BA:2247:A:H2'	23:BA:2248:C:C6	2.50	0.47
43:BU:13:VAL:HG13	43:BU:73:ARG:O	2.14	0.47
43:BU:88:LYS:HE2	43:BU:93:GLY:CA	2.39	0.47
46:DX:11:ARG:HD2	46:DX:60:PHE:HD2	1.79	0.47
23:DA:270(L):C:H6	23:DA:270(L):C:O5'	1.98	0.47
38:BP:88:ILE:HG13	38:BP:89:VAL:N	2.30	0.47
23:BA:245:G:H2'	23:BA:246:C:C6	2.49	0.47
25:BC:33:LEU:O	25:BC:35:LYS:N	2.48	0.47
23:BA:2026:C:C2	23:BA:2027:G:C8	3.03	0.47
26:BD:103:ASP:OD1	26:BD:201:THR:HG23	2.14	0.47
26:BD:169:ASN:ND2	26:BD:201:THR:HG21	2.30	0.47
34:BL:140:ALA:O	34:BL:141:ALA:CB	2.63	0.47
36:DN:10:LEU:HB2	36:DN:17:ARG:NH2	2.30	0.47
30:DH:79:ILE:HG22	30:DH:81:VAL:CG2	2.45	0.47
16:CP:27:LYS:HD2	16:CP:27:LYS:N	2.28	0.47
4:CD:8:VAL:C	4:CD:10:ARG:H	2.18	0.47
23:DA:322:A:P	27:DE:169:ASN:HB2	2.55	0.47
23:BA:2261:C:H3'	45:BW:16:SER:CB	2.45	0.47
28:DF:6:ALA:HB1	28:DF:10:LYS:HE3	1.97	0.47
4:AD:105:VAL:HG13	4:AD:110:PHE:HB2	1.96	0.47
1:AA:1188:A:H4'	14:AN:58:LYS:NZ	2.29	0.47
1:AA:9:G:H2'	1:AA:10:A:H8	1.80	0.47
18:CR:59:SER:HB3	18:CR:62:GLU:HG3	1.96	0.47
1:CA:859:A:H2'	1:CA:860:A:O4'	2.14	0.47
27:BE:157:VAL:HB	27:BE:194:MET:HB3	1.96	0.47
1:AA:859:A:H2'	1:AA:860:A:O4'	2.14	0.47
33:BK:79:PHE:HD2	38:BP:72:VAL:HG22	1.78	0.47
23:BA:636:G:OP1	34:BL:132:LYS:HD3	2.15	0.47
23:DA:2730:C:O2'	23:DA:2731:G:H5'	2.14	0.47
1:CA:731:G:H5'	1:CA:766:A:H4'	1.96	0.47
10:CJ:80:LYS:HB2	10:CJ:80:LYS:NZ	2.29	0.47
52:B4:24:THR:HG23	52:B4:27:GLY:HA3	1.96	0.47
27:DE:29:ASN:H	27:DE:112:MET:HE1	1.79	0.47
4:CD:92:VAL:O	4:CD:96:LEU:HD23	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BA:2887:U:H2'	23:BA:2888:C:C6	2.50	0.47
23:BA:2335:A:C8	23:BA:2337:G:C5	3.02	0.47
41:DS:20:VAL:HG11	41:DS:44:ALA:HA	1.97	0.47
23:BA:915:C:H2'	23:BA:916:G:C8	2.50	0.47
23:DA:433:C:H2'	23:DA:434:U:C6	2.50	0.47
28:BF:38:VAL:HG22	28:BF:93:THR:HG23	1.97	0.47
45:BW:12:ASN:O	45:BW:14:ARG:HB2	2.15	0.47
10:CJ:3:LYS:HD2	10:CJ:77:PRO:HD3	1.96	0.47
1:CA:1480:G:C5	1:CA:1481:U:C5	3.02	0.47
36:BN:31:HIS:HB2	36:BN:34:ILE:HD11	1.97	0.47
40:BR:14:VAL:HG13	40:BR:96:ILE:HG13	1.97	0.47
23:BA:329:G:H1	43:BU:19:LYS:HE3	1.79	0.47
23:DA:2190:G:H8	23:DA:2190:G:H5'	1.79	0.47
5:AE:96:PRO:HA	5:AE:117:ASP:OD2	2.15	0.47
20:AT:10:LEU:O	20:AT:13:LEU:HD13	2.14	0.47
50:B2:3:LYS:N	50:B2:3:LYS:HD2	2.29	0.47
40:DR:6:LYS:HA	40:DR:11:GLN:HB3	1.97	0.47
5:CE:76:ILE:HG12	5:CE:77:PRO:CD	2.31	0.47
46:BX:11:ARG:HD2	46:BX:60:PHE:HD2	1.79	0.47
25:DC:33:LEU:O	25:DC:35:LYS:N	2.48	0.47
9:CI:73:GLN:O	9:CI:77:ILE:HG13	2.15	0.47
23:DA:125:G:H5'	52:D4:19:ARG:HG3	1.96	0.47
25:BC:271:ILE:O	25:BC:272:ALA:CB	2.63	0.47
2:CB:28:PHE:CD2	2:CB:194:PRO:HG3	2.49	0.47
26:DD:5:LEU:C	26:DD:51:PHE:HE2	2.18	0.47
2:AB:84:GLU:HG3	2:AB:215:LEU:HB3	1.97	0.47
40:BR:78:LYS:HG3	40:BR:79:VAL:HG23	1.97	0.47
37:DO:90:GLY:O	37:DO:92:TYR:CD1	2.68	0.47
1:AA:624:C:H4'	16:AP:11:SER:N	2.29	0.47
23:DA:2850:A:H2'	23:DA:2851:A:C8	2.50	0.47
23:DA:966:G:C6	23:DA:967:C:N4	2.82	0.47
28:BF:81:LYS:C	28:BF:82:LEU:HD23	2.35	0.47
40:DR:24:LYS:HA	40:DR:92:THR:HG23	1.96	0.47
27:DE:37:VAL:HG13	27:DE:184:TYR:HD1	1.80	0.47
23:BA:1655:A:H1'	26:BD:113:PHE:CE2	2.50	0.47
1:CA:950:U:H2'	1:CA:951:G:C8	2.50	0.47
4:AD:96:LEU:HD12	4:AD:139:ARG:CD	2.45	0.47
23:DA:649:G:C5	23:DA:650:C:C4	3.03	0.47
4:CD:68:TYR:CE2	4:CD:97:LEU:HB3	2.50	0.47
11:AK:67:ASP:OD1	11:AK:71:LYS:HE3	2.15	0.47
23:DA:244:A:C2	23:DA:255:A:C4	3.03	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:857:C:H2'	1:AA:858:G:O4'	2.15	0.47
23:BA:96:G:H4'	47:BY:48:HIS:CD2	2.50	0.47
1:CA:562:C:N4	1:CA:884:U:C6	2.83	0.47
23:BA:1411:C:H2'	23:BA:1412:A:C8	2.50	0.47
1:AA:1322:C:OP2	13:AM:100:GLY:HA3	2.15	0.47
23:BA:196:A:H2'	23:BA:196:A:N3	2.29	0.47
23:BA:396:G:H1'	46:BX:42:GLN:OE1	2.15	0.47
50:D2:4:HIS:HB2	50:D2:5:PRO:HD3	1.97	0.47
34:DL:38:GLN:HG3	34:DL:41:ARG:HG2	1.96	0.47
1:AA:1239:A:H4'	1:AA:1240:U:H5'	1.96	0.47
1:CA:1239:A:H4'	1:CA:1240:U:H5'	1.97	0.47
1:CA:522:C:O2'	1:CA:523:A:H5'	2.14	0.47
34:DL:62:LEU:HA	34:DL:63:PRO:HD3	1.78	0.47
43:BU:76:CYS:O	43:BU:77:PRO:C	2.52	0.47
25:DC:136:ILE:O	25:DC:168:ARG:NH2	2.48	0.47
34:DL:61:ARG:HH11	53:D5:13:ARG:HD2	1.80	0.47
1:CA:1373:G:H5''	7:CG:36:LYS:NZ	2.29	0.47
9:CI:14:VAL:O	9:CI:65:VAL:HG23	2.14	0.47
3:CC:195:VAL:CG1	3:CC:196:LEU:N	2.77	0.47
5:CE:43:LEU:CD1	5:CE:132:ALA:HB1	2.41	0.47
2:CB:187:LEU:HD23	2:CB:201:ILE:HG22	1.96	0.47
2:AB:20:GLU:HG3	2:AB:191:ASP:H	1.80	0.47
43:BU:10:GLY:HA2	43:BU:27:VAL:HG23	1.96	0.47
17:AQ:69:LYS:O	17:AQ:70:ARG:HD2	2.15	0.47
46:BX:19:GLN:HA	46:BX:41:ARG:HA	1.96	0.47
25:DC:43:ARG:HB2	25:DC:49:ILE:HA	1.97	0.47
35:DM:35:VAL:HA	35:DM:101:ARG:O	2.14	0.47
24:BB:106:G:C6	24:BB:107:U:C4	3.03	0.47
26:BD:86:PRO:HB2	26:BD:87:GLU:H	1.41	0.47
6:AF:61:LEU:HD23	6:AF:63:TYR:OH	2.14	0.47
29:DG:123:PHE:HA	29:DG:133:VAL:HA	1.97	0.47
43:DU:76:CYS:O	43:DU:77:PRO:C	2.53	0.47
1:AA:683:G:C6	1:AA:684:A:C5	3.03	0.47
1:AA:688:G:H2'	1:AA:689:C:H6	1.80	0.47
33:BK:79:PHE:CD2	38:BP:72:VAL:HG22	2.50	0.47
24:DB:8:U:H5''	37:DO:15:ARG:NH2	2.29	0.47
13:AM:15:VAL:HG13	13:AM:43:THR:O	2.14	0.47
1:AA:731:G:H5'	1:AA:766:A:H4'	1.96	0.47
23:DA:2730:C:H4'	26:DD:168:MET:O	2.14	0.47
1:CA:963:G:H2'	1:CA:964:A:H8	1.80	0.47
23:BA:753:C:H2'	23:BA:754:C:H6	1.80	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
40:DR:99:ILE:H	40:DR:99:ILE:HD13	1.80	0.47
45:DW:42:GLY:HA2	45:DW:57:PHE:CD2	2.50	0.47
23:BA:909:A:H2'	23:BA:912:C:H5	1.79	0.47
17:AQ:85:VAL:O	17:AQ:89:LEU:HG	2.15	0.47
1:AA:67:C:H2'	1:AA:68:G:C8	2.50	0.47
1:CA:544:G:H2'	1:CA:545:C:C6	2.50	0.47
23:DA:2836:U:H2'	23:DA:2837:G:C8	2.49	0.47
23:DA:2836:U:C4	23:DA:2883:A:N6	2.83	0.47
47:DY:52:ASP:O	47:DY:56:GLN:HB2	2.14	0.47
23:DA:536:A:H2'	23:DA:537:C:C6	2.50	0.47
27:DE:132:VAL:HG23	27:DE:133:ASN:N	2.30	0.47
23:BA:1506:C:H2'	23:BA:1508:A:C8	2.50	0.47
13:AM:52:GLU:HA	13:AM:55:ARG:HB3	1.97	0.47
17:CQ:85:VAL:O	17:CQ:89:LEU:HG	2.15	0.47
42:DT:66:LEU:HD23	42:DT:67:GLY:N	2.29	0.47
11:AK:12:ARG:HG2	11:AK:13:GLN:N	2.30	0.47
36:BN:97:VAL:HA	36:BN:113:LEU:O	2.15	0.47
45:DW:32:ARG:CB	45:DW:35:ASN:HD21	2.28	0.46
7:AG:106:GLN:O	7:AG:110:GLN:HG3	2.14	0.46
34:DL:57:THR:CG2	34:DL:59:LEU:HB2	2.45	0.46
28:DF:109:VAL:C	28:DF:112:PRO:HD2	2.35	0.46
25:BC:25:THR:O	25:BC:25:THR:HG23	2.14	0.46
25:BC:25:THR:O	25:BC:27:THR:N	2.49	0.46
53:B5:54:GLU:O	53:B5:58:ILE:HG12	2.14	0.46
23:DA:773:U:H4'	25:DC:47:GLY:CA	2.40	0.46
3:CC:11:ARG:HH21	3:CC:180:ALA:HB3	1.80	0.46
2:CB:77:ALA:HB1	2:CB:165:VAL:HG11	1.96	0.46
7:CG:69:VAL:HG12	7:CG:69:VAL:O	2.15	0.46
1:AA:986:A:H1'	19:AS:54:GLY:O	2.15	0.46
33:BK:77:ILE:HD13	33:BK:78:ARG:N	2.30	0.46
30:DH:82:ARG:HD2	30:DH:89:TYR:CD2	2.50	0.46
25:DC:147:LEU:HD13	25:DC:155:LEU:CD1	2.43	0.46
25:DC:154:LYS:C	25:DC:155:LEU:HD12	2.35	0.46
17:CQ:66:SER:OG	17:CQ:69:LYS:HB3	2.14	0.46
23:DA:661:C:H4'	34:DL:16:ARG:HD2	1.96	0.46
35:BM:35:VAL:HA	35:BM:101:ARG:O	2.14	0.46
1:AA:1073:U:H2'	1:AA:1074:G:C8	2.46	0.46
1:CA:1187:G:H5'	9:CI:113:LYS:HE2	1.97	0.46
23:DA:2272:U:H5''	23:DA:2273:A:OP1	2.15	0.46
20:AT:26:ASN:HB2	20:AT:71:THR:CG2	2.44	0.46
38:BP:50:ILE:HA	38:BP:99:LEU:CD1	2.45	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1319:A:OP2	19:CS:5:LEU:HD23	2.15	0.46
38:DP:30:VAL:HG12	38:DP:86:ILE:CG1	2.45	0.46
23:DA:1952:A:C4	33:DK:22:ILE:HD12	2.49	0.46
1:CA:262:A:C6	1:CA:263:A:C6	3.03	0.46
23:BA:966:G:C4	23:BA:967:C:C5	3.03	0.46
23:DA:636:G:OP1	34:DL:132:LYS:HD3	2.15	0.46
14:AN:37:PHE:HZ	14:AN:56:VAL:HG21	1.80	0.46
1:CA:909:A:H3'	1:CA:910:C:H6	1.80	0.46
36:BN:53:HIS:O	36:BN:56:LYS:HB3	2.15	0.46
2:AB:141:GLU:O	2:AB:145:LEU:HD23	2.14	0.46
26:DD:49:LEU:O	26:DD:78:LEU:HA	2.14	0.46
5:CE:144:THR:O	5:CE:148:VAL:HG23	2.15	0.46
23:BA:532:A:C8	23:BA:2021:C:C5	3.03	0.46
6:AF:82:ARG:HA	6:AF:82:ARG:HH11	1.79	0.46
28:DF:174:GLU:HG2	28:DF:180:PHE:HD1	1.81	0.46
28:DF:38:VAL:HG22	28:DF:93:THR:HG23	1.97	0.46
1:CA:171:A:H2'	1:CA:172:A:C8	2.50	0.46
1:CA:1084:G:H5'	1:CA:1102:A:OP2	2.15	0.46
26:DD:175:VAL:O	26:DD:177:PRO:HD3	2.15	0.46
1:CA:46:G:O2'	1:CA:365:U:H1'	2.16	0.46
14:AN:6:LEU:HD22	14:AN:21:TYR:OH	2.16	0.46
25:BC:37:LEU:HD12	25:BC:38:LYS:H	1.79	0.46
24:BB:56:G:H4'	24:BB:57:A:C8	2.50	0.46
2:CB:221:LEU:O	2:CB:221:LEU:HD13	2.16	0.46
23:BA:1762:A:O5'	23:BA:1762:A:H8	1.98	0.46
1:CA:186(B):C:O2'	20:CT:89:ARG:HD2	2.15	0.46
40:DR:22:VAL:CG1	40:DR:23:GLU:N	2.76	0.46
23:BA:1019:U:H3	23:BA:114(B):A:N6	2.09	0.46
23:BA:2393:A:H5''	34:BL:62:LEU:HD12	1.97	0.46
23:DA:71:A:C2	42:DT:31:HIS:CE1	3.03	0.46
1:CA:92:G:H2'	1:CA:93:U:O4'	2.15	0.46
35:BM:75:THR:CA	35:BM:88:GLY:HA2	2.40	0.46
25:DC:25:THR:HG21	25:DC:81:ALA:HA	1.96	0.46
23:DA:2027:G:H2'	23:DA:2028:U:O4'	2.15	0.46
12:AL:82:VAL:HG21	12:AL:99:ILE:CD1	2.45	0.46
12:CL:30:PRO:HB2	12:CL:31:PHE:CD1	2.50	0.46
30:BH:12:LEU:N	30:BH:12:LEU:HD22	2.30	0.46
36:BN:10:LEU:HD23	36:BN:21:TYR:OH	2.15	0.46
23:DA:661:C:H4'	34:DL:18:ARG:HG2	1.96	0.46
1:CA:1298:C:C5	7:CG:114:ARG:NH1	2.84	0.46
23:BA:322:A:OP2	27:BE:169:ASN:HB2	2.14	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:AI:29:ASN:OD1	9:AI:64:THR:HA	2.15	0.46
9:AI:73:GLN:O	9:AI:77:ILE:HG13	2.15	0.46
23:DA:242:G:N7	53:D5:5:LYS:HG2	2.30	0.46
1:CA:32:A:C6	1:CA:33:A:C6	3.03	0.46
1:CA:506:G:C6	1:CA:507:C:C4	3.03	0.46
1:CA:1501:C:C2	1:CA:1504:G:O6	2.68	0.46
23:BA:1602:U:H3'	23:BA:1603:A:H5''	1.97	0.46
23:BA:226:G:N2	23:BA:228:A:H62	2.13	0.46
23:BA:1567:A:H5''	25:BC:58:HIS:CD2	2.50	0.46
8:AH:102:ARG:H	8:AH:102:ARG:HE	1.62	0.46
1:CA:376:G:H2'	1:CA:377:G:H8	1.80	0.46
38:DP:29:ARG:HA	38:DP:45:PHE:O	2.15	0.46
24:BB:28:C:H2'	24:BB:29:A:O4'	2.15	0.46
40:BR:34:GLU:HG3	40:BR:58:VAL:HG22	1.97	0.46
40:DR:20:LEU:HD12	40:DR:21:ARG:H	1.81	0.46
23:BA:2531:A:H2	23:BA:2658:C:O2	1.98	0.46
32:BJ:58:ARG:C	32:BJ:60:LYS:H	2.18	0.46
23:BA:1668:A:H4'	23:BA:1669:A:O5'	2.16	0.46
33:DK:76:ALA:HB3	38:DP:75:ILE:HB	1.97	0.46
23:DA:532:A:C8	23:DA:2021:C:C5	3.04	0.46
32:BJ:34:PRO:HB3	32:BJ:74:PHE:CE1	2.50	0.46
31:DI:56:ASN:HA	31:DI:59:ILE:HD12	1.98	0.46
13:AM:81:LEU:HD22	13:AM:86:CYS:SG	2.54	0.46
47:BY:31:GLU:O	47:BY:35:LEU:HB2	2.15	0.46
1:AA:515:G:H2'	1:AA:516:U:O4'	2.16	0.46
7:AG:101:LEU:O	7:AG:105:VAL:HG23	2.15	0.46
23:BA:943:U:OP1	34:BL:38:GLN:HB3	2.16	0.46
34:BL:128:HIS:CA	34:BL:147:LEU:HB3	2.22	0.46
53:D5:60:LEU:C	53:D5:62:LEU:H	2.18	0.46
9:AI:19:LEU:HD23	9:AI:20:ARG:H	1.81	0.46
10:CJ:49:VAL:HG21	14:CN:41:ARG:CB	2.45	0.46
46:DX:11:ARG:HB2	46:DX:13:ILE:CG2	2.46	0.46
25:DC:35:LYS:HD2	25:DC:35:LYS:HA	1.62	0.46
1:CA:675:A:H2'	1:CA:676:A:C8	2.51	0.46
2:AB:77:ALA:HB1	2:AB:165:VAL:HG11	1.97	0.46
1:CA:1327:C:H2'	1:CA:1328:C:H6	1.79	0.46
26:BD:201:THR:O	26:BD:202:LYS:HD3	2.15	0.46
30:DH:83:ALA:CB	30:DH:123:LEU:HD12	2.45	0.46
23:BA:2287:A:C6	23:BA:2289:G:C4	3.03	0.46
25:DC:76:PRO:CB	25:DC:116:GLN:HE21	2.28	0.46
37:BO:39:ILE:O	37:BO:48:LEU:HD13	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1295:G:H2'	1:AA:1296:C:O4'	2.15	0.46
10:AJ:55:LYS:O	10:AJ:56:HIS:CG	2.68	0.46
2:CB:168:THR:HG1	2:CB:192:SER:HA	1.81	0.46
23:BA:510:C:OP1	23:BA:511:U:OP2	2.34	0.46
23:BA:2687:U:C4	23:BA:2688:U:H5	2.33	0.46
23:BA:2666:C:H3'	23:BA:2667:C:C6	2.50	0.46
11:AK:120:ARG:HA	11:AK:121:PRO:HD3	1.79	0.46
11:CK:59:TYR:CZ	11:CK:63:LEU:HD11	2.50	0.46
23:DA:448:U:H1'	27:DE:84:VAL:CG2	2.45	0.46
27:DE:83:PHE:O	27:DE:86:GLY:N	2.44	0.46
1:AA:505:G:H2'	1:AA:506:G:C8	2.48	0.46
1:CA:1502:A:C8	1:CA:1505:G:N2	2.83	0.46
30:BH:58:LEU:HD23	30:BH:61:ARG:HD2	1.97	0.46
1:AA:793:U:H3'	1:AA:794:A:H5''	1.97	0.46
23:BA:1187:G:O5'	23:BA:1187:G:H8	1.98	0.46
1:CA:515:G:H2'	1:CA:516:U:O4'	2.15	0.46
8:CH:102:ARG:H	8:CH:102:ARG:HE	1.63	0.46
23:DA:2745:C:C4	23:DA:2746:U:C4	3.03	0.46
1:AA:818:G:C3'	1:AA:819:A:H5''	2.45	0.46
23:BA:380:U:O2'	46:BX:20:ARG:HG2	2.16	0.46
18:AR:84:LYS:HZ3	18:AR:84:LYS:HA	1.79	0.46
23:BA:2730:C:H4'	26:BD:168:MET:O	2.15	0.46
1:AA:1423:G:H2'	1:AA:1424:C:H6	1.80	0.46
23:BA:2795:G:H3'	23:BA:2797:U:H5''	1.95	0.46
23:DA:380:U:O2'	46:DX:20:ARG:HG2	2.15	0.46
23:BA:1010:A:H1'	23:BA:1153:C:C1'	2.45	0.46
30:DH:31:LEU:HD13	30:DH:37:VAL:HA	1.98	0.46
23:DA:2590:A:O2'	23:DA:2591:C:H5'	2.15	0.46
25:BC:76:PRO:CB	25:BC:116:GLN:HE21	2.28	0.46
23:DA:2887:U:H2'	23:DA:2888:C:C6	2.50	0.46
23:DA:2627:G:H8	23:DA:2627:G:O5'	1.98	0.46
26:BD:175:VAL:O	26:BD:177:PRO:HD3	2.16	0.46
27:BE:132:VAL:HG23	27:BE:133:ASN:N	2.30	0.46
33:BK:76:ALA:HB3	38:BP:75:ILE:HB	1.97	0.46
8:CH:38:ILE:HD12	8:CH:118:VAL:HG12	1.97	0.46
40:DR:75:PHE:C	40:DR:75:PHE:CD1	2.89	0.46
4:CD:131:ARG:N	4:CD:131:ARG:HD3	2.31	0.46
1:AA:619:U:C2	4:AD:135:LEU:HD21	2.50	0.46
23:DA:110:G:C2	23:DA:111:A:C8	3.03	0.46
1:CA:972:C:H4'	10:CJ:57:LYS:HG3	1.97	0.46
34:BL:128:HIS:HB3	34:BL:147:LEU:HD23	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:DL:80:TYR:CE1	34:DL:111:ARG:HG2	2.50	0.46
1:CA:520:A:N1	1:CA:536:C:H1'	2.30	0.46
2:AB:162:ILE:HD11	2:AB:184:VAL:HG13	1.97	0.46
25:DC:31:LYS:HA	25:DC:31:LYS:HD2	1.67	0.46
23:DA:2543:G:C8	23:DA:2543:G:H5'	2.43	0.46
28:DF:53:LEU:CD1	28:DF:88:ILE:HG12	2.46	0.46
23:BA:2469:A:H5'	23:BA:2470:G:OP2	2.15	0.46
23:DA:307:G:N1	23:DA:310:A:OP2	2.49	0.46
19:CS:29:ARG:HD2	19:CS:30:LEU:N	2.30	0.46
23:BA:1431:U:H2'	23:BA:1432:C:H6	1.81	0.46
28:BF:15:VAL:O	28:BF:19:LEU:HG	2.15	0.46
28:BF:6:ALA:HB1	28:BF:10:LYS:HE3	1.97	0.46
46:DX:27:GLU:HG3	46:DX:33:LYS:CD	2.45	0.46
1:CA:1151:A:O2'	1:CA:1152:A:C8	2.64	0.46
8:CH:48:TYR:HA	8:CH:60:ARG:O	2.16	0.46
44:DV:22:GLY:O	44:DV:41:LEU:HG	2.15	0.46
23:DA:2599:G:N7	25:DC:237:GLU:HG3	2.31	0.46
1:AA:625:G:H2'	1:AA:626:U:H6	1.80	0.46
23:DA:1588:C:H2'	23:DA:1589:C:C6	2.50	0.46
1:CA:1097:C:H2'	1:CA:1098:C:H6	1.80	0.46
1:CA:623:C:C4	1:CA:624:C:C5	3.04	0.46
23:DA:2320:A:C8	23:DA:2333:A:N6	2.83	0.46
21:AU:6:ARG:HG3	21:AU:15:ARG:NH1	2.30	0.46
1:CA:1413:A:C6	1:CA:1414:U:C4	3.02	0.46
23:DA:2846:G:H2'	23:DA:2847:U:H6	1.81	0.46
23:DA:69:C:H2'	23:DA:70:G:H8	1.81	0.46
1:CA:1480:G:C6	1:CA:1481:U:C4	3.04	0.46
44:DV:54:HIS:HB3	44:DV:101:PRO:HD3	1.98	0.46
1:AA:1463:C:H2'	1:AA:1464:G:H8	1.81	0.46
23:DA:2228:G:C6	23:DA:2229:C:C4	3.03	0.46
23:BA:289:A:H2'	23:BA:290:G:O4'	2.15	0.46
2:AB:52:GLU:O	2:AB:56:ARG:HG3	2.14	0.46
1:AA:579:G:H2'	1:AA:580:U:C6	2.51	0.46
1:CA:1463:C:H2'	1:CA:1464:G:H8	1.80	0.46
27:BE:167:ALA:HB1	27:BE:173:VAL:HG11	1.97	0.46
32:DJ:34:PRO:HB3	32:DJ:74:PHE:CE1	2.50	0.46
23:DA:394:A:O2'	23:DA:395:U:H5'	2.15	0.46
1:CA:110:C:H2'	1:CA:111:G:O4'	2.14	0.46
23:BA:2190:G:H5'	23:BA:2190:G:H8	1.79	0.46
23:BA:1936:A:OP1	23:BA:1936:A:H3'	2.15	0.46
16:AP:45:THR:HB	16:AP:46:PRO:HD2	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BA:1620:G:O2'	52:B4:2:LYS:HG2	2.16	0.46
23:BA:2734:A:C8	23:BA:2735:G:C8	3.04	0.46
23:BA:2056:G:C2'	23:BA:2056:G:N3	2.78	0.46
39:BQ:69:CYS:SG	39:BQ:79:PHE:HD2	2.38	0.46
32:DJ:92:GLN:O	32:DJ:94:ILE:HG13	2.15	0.46
28:DF:64:THR:HG23	28:DF:66:GLN:N	2.31	0.46
34:DL:84:ASN:HA	34:DL:115:LEU:O	2.16	0.46
47:DY:1:MET:SD	47:DY:1:MET:O	2.73	0.46
34:BL:85:LEU:HA	34:BL:88:LEU:CB	2.45	0.46
23:DA:245:G:H2'	23:DA:246:C:C6	2.44	0.46
38:BP:27:THR:HA	38:BP:48:ILE:HA	1.96	0.46
1:CA:674:G:H2'	1:CA:675:A:C8	2.47	0.46
48:BZ:28:LEU:HA	48:BZ:33:GLN:OE1	2.16	0.46
3:AC:21:ARG:O	3:AC:58:GLU:HA	2.16	0.46
23:BA:2037:G:H2'	23:BA:2038:G:C8	2.51	0.46
33:BK:71:ARG:HH21	33:BK:77:ILE:HG21	1.79	0.46
34:DL:16:ARG:CZ	34:DL:18:ARG:HG3	2.45	0.46
28:DF:15:VAL:O	28:DF:19:LEU:HG	2.16	0.46
1:AA:522:C:N4	1:AA:528:C:H42	2.14	0.46
9:AI:17:VAL:HG21	9:AI:80:GLY:C	2.36	0.46
19:AS:40:ILE:HG21	19:AS:62:ILE:HD11	1.97	0.46
23:BA:448:U:H1'	27:BE:84:VAL:CG2	2.45	0.46
18:AR:59:SER:HB3	18:AR:62:GLU:HG3	1.96	0.46
47:DY:41:ILE:HD11	47:DY:44:LEU:HD12	1.96	0.46
27:BE:124:LEU:HD12	27:BE:125:LEU:N	2.31	0.46
33:BK:31:LYS:HB3	33:BK:32:TYR:CE1	2.51	0.46
1:AA:192:U:O2'	1:AA:193:C:H5'	2.16	0.46
1:CA:1242:C:O2'	1:CA:1303:C:H5''	2.15	0.46
23:BA:2830:G:N3	23:BA:2883:A:H2	2.13	0.46
8:AH:23:SER:HB3	8:AH:62:TYR:HA	1.97	0.46
36:DN:33:ARG:HG3	36:DN:115:GLU:HG3	1.98	0.46
32:BJ:105:LEU:O	32:BJ:106:LYS:C	2.54	0.46
27:DE:54:ARG:HA	27:DE:87:GLY:HA3	1.97	0.46
1:CA:767:A:H2'	1:CA:768:A:O4'	2.15	0.46
1:AA:1015:A:H8	1:AA:1015:A:O5'	1.98	0.46
45:BW:56:ASP:O	45:BW:57:PHE:HB2	2.16	0.46
1:AA:1333:A:H2'	1:AA:1334:G:O4'	2.16	0.46
23:BA:587:C:H42	34:BL:33:ARG:CG	2.18	0.46
5:AE:77:PRO:HG2	5:AE:78:HIS:ND1	2.30	0.46
42:DT:26:TYR:HE1	42:DT:83:VAL:HG21	1.80	0.46
23:DA:1019:U:H3	23:DA:114(B):A:N6	2.10	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BA:958:U:H5'	35:BM:14:ARG:NH1	2.30	0.46
44:BV:74:VAL:O	44:BV:76:LEU:HD12	2.16	0.46
23:DA:1173:G:H3'	23:DA:1174:A:H5''	1.97	0.46
23:BA:1173:G:H3'	23:BA:1174:A:H5''	1.96	0.46
28:BF:86:MET:O	28:BF:87:PRO:O	2.34	0.46
28:BF:53:LEU:CD1	28:BF:88:ILE:HG12	2.46	0.46
30:DH:83:ALA:CA	30:DH:89:TYR:HD1	2.28	0.46
36:BN:21:TYR:CE2	36:BN:43:GLU:HB3	2.50	0.46
23:DA:1448:G:H21	23:DA:1529:A:H2	1.63	0.46
46:BX:27:GLU:HG2	46:BX:28:GLY:N	2.30	0.46
1:AA:1371:G:OP1	9:AI:11:LYS:HB3	2.15	0.46
1:CA:668:G:H1'	15:CO:46:HIS:HD2	1.81	0.46
5:CE:102:ALA:HB2	5:CE:120:THR:OG1	2.16	0.46
29:DG:58:GLU:HB2	29:DG:61:HIS:ND1	2.30	0.46
9:AI:62:TYR:C	9:AI:63:ILE:HD12	2.36	0.46
30:BH:15:VAL:HG12	30:BH:16:GLY:N	2.30	0.46
29:DG:103:LEU:HD22	29:DG:123:PHE:CE1	2.50	0.46
12:AL:84:ILE:HG23	12:AL:97:TYR:HB3	1.98	0.46
23:DA:2667:C:H1'	29:DG:109:PHE:HD2	1.81	0.46
1:CA:828:A:H5''	1:CA:859:A:C2	2.50	0.46
38:BP:30:VAL:HG12	38:BP:86:ILE:CG1	2.45	0.46
23:BA:2745:C:C4	23:BA:2746:U:C4	3.03	0.46
1:AA:238:G:P	17:AQ:25:ARG:HH22	2.39	0.46
1:AA:1505:G:H4'	1:AA:1506:U:H5'	1.98	0.46
23:BA:2815:C:O2'	50:B2:42:PRO:HB2	2.16	0.46
2:CB:22:LYS:HZ3	2:CB:22:LYS:N	2.13	0.46
5:CE:12:LEU:C	5:CE:13:ILE:HD12	2.36	0.46
23:BA:69:C:H2'	23:BA:70:G:C8	2.50	0.46
13:AM:70:LEU:C	13:AM:70:LEU:HD23	2.36	0.46
23:DA:96:G:H4'	47:DY:48:HIS:CD2	2.51	0.46
44:BV:28:MET:HE2	44:BV:28:MET:HB3	1.80	0.46
27:DE:143:ALA:HB1	27:DE:148:LEU:HB2	1.97	0.46
1:CA:67:C:H2'	1:CA:68:G:C8	2.50	0.46
50:D2:18:ALA:O	50:D2:21:SER:HB2	2.16	0.46
23:DA:836:G:H2'	23:DA:837:C:C6	2.51	0.46
23:DA:1762:A:H8	23:DA:1762:A:O5'	1.97	0.46
1:AA:1442:G:H3'	1:AA:1442:G:H8	1.81	0.46
48:DZ:55:ARG:HA	48:DZ:55:ARG:HD3	1.65	0.46
23:DA:1164:G:C5'	23:DA:1164:G:C8	2.99	0.46
25:DC:132:PRO:HD3	25:DC:190:TYR:CE2	2.51	0.46
27:DE:63:LYS:HG2	27:DE:65:TRP:O	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:CB:20:GLU:HG3	2:CB:191:ASP:H	1.80	0.46
7:CG:69:VAL:O	7:CG:71:PRO:HD3	2.16	0.46
23:DA:2287:A:C6	23:DA:2289:G:C4	3.02	0.46
28:DF:86:MET:O	28:DF:87:PRO:O	2.33	0.46
25:DC:130:ALA:HA	25:DC:192:THR:HA	1.96	0.46
4:AD:8:VAL:C	4:AD:10:ARG:H	2.18	0.46
51:B3:34:LEU:N	51:B3:34:LEU:HD13	2.31	0.46
40:DR:77:ALA:C	40:DR:79:VAL:H	2.19	0.46
25:BC:69:ARG:HH12	25:BC:117:VAL:CG2	2.29	0.46
23:DA:2273:A:H2'	23:DA:2274:A:C8	2.51	0.46
23:DA:2582:G:C2	23:DA:2583:G:C8	3.03	0.46
10:CJ:6:ILE:HG22	10:CJ:98:ILE:CG2	2.45	0.46
1:CA:804:U:H5''	1:CA:805:C:OP2	2.16	0.46
23:BA:2476:A:C2	23:BA:2477:C:C6	3.03	0.46
33:DK:79:PHE:CD2	38:DP:72:VAL:HG22	2.50	0.46
1:AA:828:A:H5''	1:AA:859:A:C2	2.50	0.46
40:BR:24:LYS:HA	40:BR:92:THR:HG23	1.97	0.46
39:DQ:28:ARG:HG3	39:DQ:38:THR:OG1	2.15	0.46
5:AE:11:ILE:HG21	5:AE:105:VAL:HG22	1.98	0.46
1:AA:950:U:H2'	1:AA:951:G:C8	2.50	0.46
35:DM:47:ILE:HG22	35:DM:48:GLU:N	2.30	0.46
23:DA:2012:G:O2'	41:DS:96:ILE:HD11	2.16	0.46
23:BA:2320:A:C8	23:BA:2333:A:N6	2.83	0.46
1:AA:924:C:H2'	1:AA:925:G:C8	2.50	0.46
45:DW:50:ASN:ND2	45:DW:83:PRO:HD3	2.30	0.46
26:BD:78:LEU:N	26:BD:78:LEU:HD23	2.30	0.46
1:AA:102:G:H2'	1:AA:103:C:H6	1.81	0.46
3:AC:8:ILE:CD1	3:AC:16:ARG:HH21	2.29	0.46
41:DS:9:TYR:H	41:DS:102:HIS:CD2	2.34	0.46
7:CG:45:ASP:O	7:CG:49:ILE:HG12	2.16	0.46
23:BA:1449:G:H2'	23:BA:1450:C:C6	2.50	0.46
38:BP:63:VAL:O	38:BP:73:GLU:HA	2.16	0.46
1:CA:814:A:N7	1:CA:816:A:C4	2.84	0.46
23:BA:2590:A:O2'	23:BA:2591:C:H5'	2.15	0.46
1:AA:110:C:H2'	1:AA:111:G:O4'	2.15	0.46
40:DR:34:GLU:HG3	40:DR:58:VAL:HG22	1.98	0.46
23:BA:394:A:O2'	23:BA:395:U:H5'	2.16	0.46
23:DA:286:C:H2'	23:DA:287:C:H6	1.81	0.46
1:AA:509:A:C6	1:AA:510:A:N1	2.84	0.46
1:AA:324:G:N2	1:AA:327:A:C8	2.83	0.46
23:DA:1121:C:O5'	23:DA:1121:C:H6	1.99	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
32:DJ:81:ASP:OD1	32:DJ:81:ASP:N	2.49	0.46
23:DA:1632:A:H8	23:DA:1632:A:O5'	1.99	0.46
13:CM:52:GLU:HA	13:CM:55:ARG:HB3	1.98	0.46
23:BA:613:U:H4'	23:BA:616:A:N6	2.31	0.46
23:BA:258:G:H2'	23:BA:259:G:H8	1.80	0.46
34:BL:38:GLN:CG	34:BL:39:LYS:H	2.28	0.46
42:BT:55:ASN:HB2	42:BT:80:ILE:HG23	1.97	0.46
28:DF:60:LEU:HD13	28:DF:60:LEU:C	2.36	0.46
51:D3:25:LYS:HD3	53:D5:34:TRP:CZ3	2.50	0.46
37:DO:49:VAL:CG1	37:DO:76:LYS:HB2	2.46	0.46
49:D1:57:ILE:HG22	49:D1:59:VAL:CG2	2.46	0.46
10:AJ:75:ILE:HG13	10:AJ:76:ASN:N	2.20	0.46
9:CI:77:ILE:O	9:CI:81:ILE:HG13	2.15	0.46
1:AA:321:A:C2	1:AA:333:G:C2	3.04	0.46
7:AG:69:VAL:O	7:AG:71:PRO:HD3	2.15	0.46
28:DF:88:ILE:HG13	28:DF:89:GLY:N	2.31	0.46
30:BH:78:THR:HA	30:BH:143:SER:CB	2.44	0.46
19:CS:29:ARG:O	19:CS:31:ILE:HG22	2.16	0.46
23:DA:558:G:H2'	23:DA:559:G:H8	1.80	0.46
4:CD:9:CYS:HB3	4:CD:32:ALA:CB	2.45	0.46
23:BA:1678:G:H22	23:BA:1989:G:H22	1.62	0.46
46:BX:27:GLU:HG3	46:BX:33:LYS:CD	2.46	0.46
1:AA:729:A:H2'	1:AA:730:G:C8	2.45	0.46
1:AA:1254:C:OP1	10:AJ:45:ARG:HD3	2.15	0.46
23:DA:1348:G:C2'	23:DA:1349:A:H5''	2.45	0.46
24:BB:79:C:O5'	24:BB:79:C:H6	1.99	0.46
23:BA:2273:A:H2'	23:BA:2274:A:C8	2.50	0.46
1:AA:1188:A:H2'	1:AA:1189:C:O4'	2.16	0.46
27:DE:198:ALA:O	27:DE:201:VAL:HG12	2.16	0.46
23:DA:36:G:N1	23:DA:445:C:C4	2.84	0.46
13:AM:27:LYS:HE2	13:AM:31:LYS:CE	2.45	0.46
1:CA:892:A:H2'	1:CA:893:C:H6	1.77	0.46
23:BA:221:A:N7	23:BA:266:G:C5	2.84	0.46
23:BA:534:U:O2'	39:BQ:49:HIS:CD2	2.69	0.46
14:CN:37:PHE:HZ	14:CN:56:VAL:HG21	1.81	0.46
1:CA:790:A:H5'	22:CV:6168:G:H4'	1.97	0.46
23:DA:1178:C:H2'	23:DA:1179:C:C6	2.51	0.46
1:CA:1432:G:H8	1:CA:1432:G:O5'	1.99	0.46
23:DA:1354:A:H2'	23:DA:1355:G:O4'	2.16	0.46
28:BF:39:ILE:HG12	28:BF:157:ILE:HG22	1.98	0.46
1:AA:1410:G:H2'	1:AA:1411:C:C6	2.51	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
38:BP:62:THR:HG22	38:BP:75:ILE:HG13	1.98	0.46
13:AM:23:TYR:HE1	13:AM:70:LEU:HD22	1.81	0.46
41:DS:9:TYR:H	41:DS:102:HIS:HD2	1.62	0.46
9:CI:29:ASN:OD1	9:CI:64:THR:HA	2.15	0.46
23:BA:451:C:H41	23:BA:453:C:H3'	1.80	0.46
23:BA:2764:A:N7	23:BA:2766:G:C6	2.83	0.46
7:AG:74:GLU:HG2	7:AG:91:VAL:HG22	1.98	0.46
11:CK:12:ARG:HG2	11:CK:13:GLN:N	2.30	0.46
1:AA:596:C:H6	1:AA:596:C:H5'	1.81	0.46
27:BE:50:SER:HB2	27:BE:94:PRO:HD3	1.97	0.46
1:CA:224:C:H2'	1:CA:225:C:C6	2.51	0.46
1:AA:520:A:N1	1:AA:536:C:H1'	2.31	0.46
38:BP:29:ARG:HA	38:BP:45:PHE:O	2.16	0.46
14:CN:6:LEU:HD22	14:CN:21:TYR:OH	2.16	0.46
39:DQ:69:CYS:CB	39:DQ:79:PHE:HD2	2.29	0.46
39:BQ:69:CYS:CB	39:BQ:79:PHE:HD2	2.29	0.46
42:DT:25:LYS:HE3	42:DT:82:GLN:OE1	2.16	0.46
32:BJ:92:GLN:O	32:BJ:94:ILE:HG13	2.16	0.46
23:BA:71:A:OP1	23:BA:72:U:H2'	2.16	0.46
23:DA:958:U:H5'	35:DM:14:ARG:NH1	2.31	0.46
27:BE:63:LYS:HG2	27:BE:65:TRP:O	2.15	0.46
46:BX:11:ARG:HB2	46:BX:13:ILE:CG2	2.45	0.46
17:AQ:54:GLY:HA3	17:AQ:82:MET:CE	2.46	0.46
36:DN:21:TYR:CE2	36:DN:43:GLU:HB3	2.50	0.46
23:DA:2039:C:C2	23:DA:2040:C:C5	3.04	0.46
17:CQ:45:HIS:HB2	17:CQ:69:LYS:HE2	1.98	0.46
23:BA:558:G:H2'	23:BA:559:G:H8	1.81	0.46
19:CS:49:ILE:N	19:CS:49:ILE:HD12	2.31	0.46
17:AQ:45:HIS:HB2	17:AQ:69:LYS:HE2	1.98	0.46
23:BA:661:C:H4'	34:BL:18:ARG:HG2	1.98	0.46
23:BA:1317:A:N6	23:BA:1336:A:N6	2.64	0.46
19:AS:29:ARG:HD2	19:AS:30:LEU:N	2.30	0.46
51:B3:34:LEU:HD23	51:B3:36:LEU:HD22	1.98	0.46
23:DA:478:A:C6	23:DA:480:A:C6	3.03	0.46
26:BD:118:LYS:NZ	36:BN:2:ARG:HH22	2.13	0.46
23:DA:1654:A:OP2	36:DN:3:HIS:NE2	2.49	0.46
19:CS:10:PHE:O	19:CS:11:VAL:HB	2.16	0.46
1:AA:622:A:C8	1:AA:623:C:C6	3.04	0.46
23:BA:2666:C:H3'	23:BA:2667:C:H6	1.81	0.46
23:BA:2630:G:H1'	23:BA:2894:G:H1'	1.98	0.46
1:CA:505:G:H2'	1:CA:506:G:C8	2.49	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:BC:70:TRP:CH2	25:BC:150:LYS:HA	2.51	0.46
25:BC:146:GLU:OE2	25:BC:150:LYS:N	2.49	0.46
28:DF:81:LYS:C	28:DF:82:LEU:HD23	2.37	0.46
23:BA:2647:U:H2'	23:BA:2648:C:H6	1.80	0.46
23:BA:2306:C:H4'	28:BF:136:ARG:HH22	1.80	0.46
1:CA:1191:A:H5''	3:CC:4:LYS:HZ2	1.81	0.46
23:BA:1833:U:C2	23:BA:1834:U:C5	3.04	0.46
23:BA:499:U:C4'	43:BU:47:LYS:HZ1	2.29	0.46
8:AH:35:ILE:O	8:AH:39:LEU:HB2	2.16	0.46
27:DE:28:ILE:HA	27:DE:112:MET:HE3	1.97	0.46
38:BP:109:GLU:HA	38:BP:112:ARG:HG3	1.98	0.46
23:BA:2252:G:H2'	23:BA:2253:G:C8	2.51	0.46
23:BA:155:C:H2'	23:BA:161:U:H5'	1.98	0.46
2:CB:141:GLU:O	2:CB:145:LEU:HD23	2.15	0.46
45:BW:14:ARG:O	45:BW:15:ASP:HB2	2.16	0.46
35:BM:134:ARG:O	35:BM:135:ASP:HB2	2.16	0.46
23:BA:1354:A:C8	23:BA:1355:G:C8	3.03	0.46
26:DD:190:GLY:HA2	26:DD:191:PRO:HD3	1.75	0.46
5:CE:96:PRO:HA	5:CE:117:ASP:OD2	2.16	0.46
23:BA:2678:C:H2'	23:BA:2679:A:H8	1.81	0.46
23:DA:1252:G:C2	23:DA:1253:A:C2	3.04	0.46
27:DE:70:THR:HG23	27:DE:72:ARG:H	1.81	0.46
9:CI:52:ALA:C	9:CI:95:LYS:HZ1	2.19	0.46
2:AB:73:THR:HA	2:AB:94:ASN:O	2.16	0.46
1:AA:224:C:H2'	1:AA:225:C:C6	2.51	0.46
12:AL:110:LYS:O	12:AL:111:ASP:HB2	2.16	0.46
23:DA:2527:C:H6	23:DA:2527:C:O5'	1.98	0.46
47:DY:59:ARG:HG2	47:DY:59:ARG:H	1.59	0.46
2:AB:221:LEU:O	2:AB:221:LEU:HD13	2.16	0.46
32:BJ:81:ASP:OD1	32:BJ:81:ASP:N	2.48	0.46
23:BA:2722:G:H4'	36:BN:5:LYS:HB3	1.98	0.46
23:DA:2315:G:H2'	23:DA:2316:C:C6	2.51	0.46
23:BA:587:C:C5	23:BA:671:C:H1'	2.51	0.46
23:BA:826:U:C5	23:BA:828:U:H1'	2.50	0.46
23:BA:833:U:H5''	34:BL:48:PRO:HB2	1.97	0.46
7:CG:113:GLU:HB2	7:CG:119:ARG:CG	2.28	0.46
39:DQ:69:CYS:SG	39:DQ:79:PHE:HD2	2.39	0.46
42:BT:24:GLY:O	42:BT:83:VAL:HG22	2.15	0.46
28:BF:64:THR:HG23	28:BF:66:GLN:N	2.30	0.46
1:CA:1224:G:C4'	13:CM:102:ARG:HH22	2.24	0.46
49:B1:59:VAL:HG12	49:B1:60:GLU:N	2.19	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
44:DV:71:VAL:HG11	44:DV:74:VAL:CG2	2.46	0.46
16:AP:4:ILE:HD12	16:AP:4:ILE:N	2.31	0.46
25:BC:102:LYS:C	25:BC:103:ARG:HG2	2.37	0.46
1:AA:321:A:N7	1:AA:328:C:C6	2.84	0.46
23:BA:1006:C:H5'	32:BJ:51:THR:HG23	1.97	0.46
40:BR:79:VAL:HG12	40:BR:79:VAL:O	2.16	0.46
37:BO:34:HIS:CG	37:BO:54:LEU:HB2	2.51	0.46
4:AD:9:CYS:HB3	4:AD:32:ALA:CB	2.44	0.46
4:CD:117:ALA:O	4:CD:121:VAL:HG23	2.16	0.46
25:BC:130:ALA:HA	25:BC:192:THR:HA	1.97	0.46
5:AE:91:LEU:N	5:AE:91:LEU:HD12	2.30	0.46
44:BV:24:LEU:HD21	44:BV:86:VAL:HG23	1.98	0.46
23:DA:863:A:OP1	35:DM:21:THR:HB	2.16	0.46
23:BA:2667:C:H1'	29:BG:109:PHE:HD2	1.81	0.46
4:CD:64:LEU:HB2	4:CD:198:VAL:HG11	1.98	0.46
1:CA:947:G:H2'	1:CA:948:C:C6	2.51	0.46
38:DP:20:PRO:CD	38:DP:86:ILE:HG23	2.45	0.46
23:DA:414:C:H2'	23:DA:415:A:H8	1.80	0.46
28:DF:128:ARG:HH21	28:DF:130:ASN:HD21	1.63	0.46
1:AA:892:A:H2'	1:AA:893:C:H6	1.80	0.46
23:BA:919:G:H2'	23:BA:920:G:C8	2.50	0.46
41:DS:10:VAL:O	41:DS:12:ILE:N	2.49	0.46
6:AF:62:TRP:CB	18:AR:35:ARG:HH12	2.29	0.46
2:AB:131:PRO:O	2:AB:135:GLN:HG3	2.16	0.46
2:AB:138:LEU:O	2:AB:141:GLU:HB2	2.16	0.46
4:CD:96:LEU:HD12	4:CD:139:ARG:CD	2.45	0.46
23:DA:356:G:H2'	23:DA:357:A:H8	1.81	0.46
45:DW:12:ASN:O	45:DW:14:ARG:N	2.49	0.46
1:AA:1480:G:C6	1:AA:1481:U:C4	3.04	0.46
23:DA:2830:G:N3	23:DA:2883:A:H2	2.14	0.46
44:DV:54:HIS:CG	44:DV:101:PRO:HG3	2.51	0.46
45:BW:42:GLY:HA2	45:BW:57:PHE:CD2	2.51	0.46
23:DA:239:U:O2'	23:DA:240:G:H5'	2.15	0.46
8:AH:114:THR:HG21	8:AH:119:LEU:HD12	1.98	0.46
29:DG:24:VAL:HG23	29:DG:37:VAL:HG21	1.98	0.46
1:AA:843:U:H5'	1:AA:848:C:C6	2.51	0.46
26:BD:24:THR:HB	26:BD:186:GLY:O	2.16	0.46
23:BA:2627:G:H8	23:BA:2627:G:O5'	1.99	0.46
1:CA:39:G:C2	1:CA:40:C:C6	3.04	0.46
23:DA:1487:G:H2'	23:DA:1488:G:H8	1.80	0.46
23:DA:1748:G:H2'	23:DA:1749:A:C8	2.51	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:AD:103:ASN:OD1	4:AD:114:ARG:NH2	2.49	0.46
4:AD:131:ARG:HD3	4:AD:131:ARG:N	2.31	0.46
24:DB:56:G:H4'	24:DB:57:A:C8	2.50	0.46
27:DE:167:ALA:HB1	27:DE:173:VAL:HG11	1.97	0.46
43:BU:68:HIS:ND1	43:BU:70:SER:HB3	2.31	0.46
27:BE:89:VAL:HG12	27:BE:90:PHE:CD2	2.51	0.45
40:BR:7:THR:HG23	40:BR:22:VAL:HG11	1.98	0.45
23:BA:1164:G:C8	23:BA:1164:G:C5'	2.99	0.45
47:DY:33:MET:O	47:DY:37:PHE:HB2	2.15	0.45
2:AB:162:ILE:O	2:AB:185:ILE:HG12	2.16	0.45
23:BA:2392:A:H2	23:BA:2424:C:N4	2.11	0.45
34:DL:65:ARG:HD2	34:DL:65:ARG:H	1.81	0.45
37:BO:72:ALA:O	37:BO:76:LYS:HG3	2.16	0.45
38:BP:46:GLU:OE2	38:BP:89:VAL:HG11	2.16	0.45
1:AA:81:G:C5	1:AA:82:U:C4	3.04	0.45
23:BA:2027:G:C2	23:BA:2028:U:H1'	2.51	0.45
23:BA:2033:A:O2'	23:BA:2034:U:P	2.74	0.45
43:DU:11:ASP:H	43:DU:27:VAL:CG2	2.29	0.45
23:DA:1210:A:H8	23:DA:1210:A:C5'	2.26	0.45
28:BF:25:TYR:CZ	28:BF:32:PRO:HD3	2.51	0.45
40:BR:72:VAL:CG2	40:BR:85:LYS:HB3	2.46	0.45
11:CK:87:THR:HA	11:CK:91:ARG:NH2	2.29	0.45
1:AA:404:U:H2'	1:AA:405:U:H6	1.81	0.45
4:AD:3:ARG:HD3	4:AD:5:ILE:HD13	1.97	0.45
51:D3:34:LEU:HD23	51:D3:36:LEU:HD22	1.98	0.45
1:CA:404:U:H2'	1:CA:405:U:H6	1.80	0.45
4:CD:3:ARG:HD3	4:CD:5:ILE:HD13	1.98	0.45
1:CA:1188:A:H2'	1:CA:1189:C:O4'	2.16	0.45
20:AT:71:THR:HG22	20:AT:72:LEU:N	2.32	0.45
11:CK:57:THR:HG22	11:CK:59:TYR:N	2.30	0.45
27:BE:198:ALA:O	27:BE:201:VAL:HG12	2.15	0.45
1:CA:1505:G:H4'	1:CA:1506:U:H5'	1.98	0.45
35:DM:66:ILE:HG22	35:DM:104:PHE:HD2	1.80	0.45
27:DE:11:VAL:HG13	27:DE:196:LEU:HD21	1.97	0.45
27:DE:157:VAL:HB	27:DE:194:MET:HB3	1.97	0.45
23:BA:1378:A:H4'	23:BA:1379:A:OP1	2.15	0.45
1:CA:819:A:N6	1:CA:1529:G:C5	2.84	0.45
46:DX:45:ASN:HD21	46:DX:47:GLN:NE2	2.15	0.45
25:DC:174:ILE:HD12	25:DC:174:ILE:N	2.31	0.45
23:DA:1010:A:H1'	23:DA:1153:C:C1'	2.46	0.45
45:DW:14:ARG:O	45:DW:15:ASP:HB2	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
31:BI:56:ASN:HA	31:BI:59:ILE:HD12	1.98	0.45
23:BA:1748:G:H2'	23:BA:1749:A:C8	2.51	0.45
11:AK:108:ILE:O	18:AR:87:ARG:HA	2.16	0.45
13:CM:23:TYR:HE1	13:CM:70:LEU:HD22	1.82	0.45
34:BL:65:ARG:H	34:BL:65:ARG:HD2	1.80	0.45
1:AA:705:U:C5	1:AA:706:A:C5	3.03	0.45
11:CK:67:ASP:OD1	11:CK:71:LYS:HE3	2.16	0.45
23:BA:1599:C:OP2	42:BT:36:LYS:HD3	2.15	0.45
29:BG:38:SER:HB2	29:BG:41:MET:HG3	1.98	0.45
36:DN:31:HIS:HB2	36:DN:34:ILE:HD11	1.97	0.45
23:DA:943:U:OP1	34:DL:38:GLN:HB3	2.16	0.45
23:DA:1826:G:H2'	23:DA:1827:C:C6	2.52	0.45
23:DA:1970:A:H4'	23:DA:1971:A:OP1	2.15	0.45
43:DU:13:VAL:CG1	43:DU:72:VAL:HB	2.47	0.45
43:DU:15:VAL:HG22	43:DU:72:VAL:HG12	1.98	0.45
23:DA:1022:G:N2	23:DA:114(B):A:C2	2.84	0.45
42:BT:30:VAL:HG12	42:BT:31:HIS:H	1.80	0.45
1:CA:922:G:C6	1:CA:923:A:C6	3.04	0.45
23:BA:956:G:H22	23:BA:959:A:H3'	1.81	0.45
49:B1:57:ILE:HG22	49:B1:59:VAL:CG2	2.46	0.45
43:DU:88:LYS:HE2	43:DU:93:GLY:CA	2.42	0.45
23:BA:947:G:H2'	23:BA:948:G:C5'	2.43	0.45
1:CA:1279:A:N6	3:CC:26:LYS:HE2	2.24	0.45
23:DA:1006:C:H5'	32:DJ:51:THR:HG23	1.98	0.45
23:BA:125:G:H5'	52:B4:19:ARG:HG3	1.97	0.45
34:DL:30:THR:HG22	34:DL:31:ALA:H	1.79	0.45
2:CB:84:GLU:HG3	2:CB:215:LEU:HB3	1.96	0.45
1:CA:392:G:C2	1:CA:393:A:C4	3.04	0.45
25:DC:228:PRO:HD3	25:DC:234:GLY:O	2.16	0.45
25:DC:76:PRO:HB3	25:DC:116:GLN:HE21	1.81	0.45
1:AA:522:C:O2'	1:AA:523:A:H5'	2.16	0.45
1:AA:1372:U:OP1	9:AI:71:SER:HB3	2.16	0.45
23:DA:2359:C:H2'	23:DA:2360:A:H8	1.81	0.45
23:BA:2582:G:C2	23:BA:2583:G:C8	3.03	0.45
25:DC:5:LYS:H	25:DC:5:LYS:HD2	1.80	0.45
1:CA:1245:A:OP2	21:CU:9:ARG:NH2	2.49	0.45
32:BJ:62:ARG:CZ	32:BJ:64:ASP:HB2	2.46	0.45
1:AA:1319:A:OP2	19:AS:5:LEU:HD23	2.15	0.45
23:BA:414:C:H2'	23:BA:415:A:H8	1.79	0.45
44:BV:58:VAL:HG11	44:BV:66:SER:HB2	1.98	0.45
1:CA:1296:C:H5''	1:CA:1297:C:OP2	2.15	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DA:581:C:H2'	23:DA:582:G:C8	2.51	0.45
1:AA:99:C:C6	1:AA:99:C:C3'	3.00	0.45
1:CA:625:G:H2'	1:CA:626:U:H6	1.81	0.45
23:BA:270(Q):C:O2'	23:BA:270(R):C:H6	1.98	0.45
1:CA:270:A:C6	1:CA:271:C:C4	3.04	0.45
23:BA:628:G:H2'	23:BA:629:G:C8	2.51	0.45
1:AA:1238:A:C8	1:AA:1303:C:H1'	2.51	0.45
21:AU:14:TRP:CE3	21:AU:15:ARG:HG2	2.51	0.45
44:BV:10:ARG:HG2	44:BV:11:GLU:N	2.32	0.45
23:DA:357:A:H2'	23:DA:358:U:C6	2.51	0.45
41:BS:9:TYR:H	41:BS:102:HIS:CD2	2.34	0.45
23:BA:1750:G:H2'	23:BA:1751:C:C6	2.52	0.45
23:DA:532:A:N1	23:DA:2020:A:H1'	2.30	0.45
13:CM:70:LEU:C	13:CM:70:LEU:HD23	2.36	0.45
1:AA:39:G:C2	1:AA:40:C:C6	3.04	0.45
1:CA:1333:A:H2'	1:CA:1334:G:O4'	2.15	0.45
1:AA:337:C:H2'	1:AA:338:A:H8	1.80	0.45
1:AA:376:G:H2'	1:AA:377:G:H8	1.80	0.45
23:DA:2722:G:H4'	36:DN:5:LYS:HB3	1.97	0.45
1:AA:384:G:H2'	1:AA:385:C:C6	2.51	0.45
23:BA:2585:U:H4'	23:BA:2586:C:OP1	2.16	0.45
23:DA:2655:G:N2	23:DA:2664:G:C5	2.84	0.45
32:BJ:63:PRO:O	39:BQ:64:ARG:HD2	2.17	0.45
23:DA:1270:C:H5''	23:DA:1271:G:O5'	2.16	0.45
23:BA:1850:G:C6	23:BA:1851:U:C4	3.04	0.45
1:CA:775:G:H2'	1:CA:776:G:O4'	2.16	0.45
28:BF:174:GLU:HG2	28:BF:180:PHE:CD1	2.51	0.45
53:B5:29:LYS:NZ	53:B5:29:LYS:HB3	2.31	0.45
33:BK:99:PHE:CD1	33:BK:99:PHE:N	2.85	0.45
12:AL:78:GLU:HG2	12:AL:78:GLU:O	2.16	0.45
24:DB:6:C:C2	24:DB:115:G:N2	2.84	0.45
15:AO:62:GLN:O	15:AO:66:LEU:HD13	2.16	0.45
26:BD:54:GLN:HB2	26:BD:74:PRO:O	2.16	0.45
1:AA:1448:C:H2'	1:AA:1449:C:H6	1.81	0.45
23:DA:868:U:C4	23:DA:869:G:N7	2.85	0.45
34:BL:50:ARG:HD3	34:BL:51:PHE:HB2	1.98	0.45
9:CI:19:LEU:HD23	9:CI:20:ARG:H	1.81	0.45
42:BT:54:VAL:C	42:BT:55:ASN:HD22	2.20	0.45
42:BT:26:TYR:HE1	42:BT:83:VAL:HG21	1.81	0.45
28:BF:60:LEU:C	28:BF:60:LEU:HD13	2.37	0.45
25:BC:132:PRO:HD3	25:BC:190:TYR:CE2	2.52	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:BC:31:LYS:HA	25:BC:31:LYS:HD2	1.65	0.45
26:BD:101:ARG:HB3	26:BD:169:ASN:HD22	1.82	0.45
23:DA:388:G:OP1	46:DX:33:LYS:HB3	2.16	0.45
11:CK:33:THR:HA	11:CK:40:ILE:HG12	1.97	0.45
40:BR:72:VAL:HG22	40:BR:85:LYS:O	2.16	0.45
4:CD:173:TRP:NE1	4:CD:189:PRO:HG3	2.31	0.45
23:DA:1331:A:O2'	23:DA:1332:G:C8	2.68	0.45
23:DA:2346:A:H5''	23:DA:2383:G:C1'	2.45	0.45
1:AA:1349:A:H2'	1:AA:1350:A:O4'	2.17	0.45
4:AD:64:LEU:HB2	4:AD:198:VAL:HG11	1.98	0.45
1:CA:793:U:H3'	1:CA:794:A:H5''	1.97	0.45
23:BA:779:U:P	25:BC:49:ILE:HG13	2.56	0.45
27:BE:192:LEU:C	27:BE:192:LEU:HD23	2.37	0.45
23:BA:1588:C:H2'	23:BA:1589:C:C6	2.49	0.45
1:CA:1295:G:H2'	1:CA:1296:C:O4'	2.16	0.45
23:DA:301:G:H5'	23:DA:334:C:O2'	2.15	0.45
1:AA:235:C:H2'	1:AA:236:G:H8	1.81	0.45
1:CA:192:U:H2'	1:CA:193:C:H6	1.82	0.45
23:BA:1995:U:H3'	23:BA:1996:C:H2'	1.99	0.45
1:AA:819:A:N6	1:AA:1529:G:C5	2.84	0.45
23:DA:919:G:C5'	24:DB:81:G:H1'	2.47	0.45
1:CA:1238:A:C8	1:CA:1303:C:H1'	2.51	0.45
3:AC:33:LEU:HD21	14:AN:53:LEU:CD2	2.46	0.45
6:CF:17:SER:O	6:CF:21:LEU:HD23	2.16	0.45
23:BA:1709:U:C2	23:BA:1750:G:C2	3.05	0.45
45:BW:12:ASN:O	45:BW:14:ARG:N	2.49	0.45
28:DF:174:GLU:HG2	28:DF:180:PHE:CD1	2.51	0.45
23:BA:1354:A:H2'	23:BA:1355:G:O4'	2.17	0.45
23:DA:969:U:H2'	23:DA:970:C:C6	2.52	0.45
23:BA:447:A:C5	23:BA:473:G:C5	3.05	0.45
41:BS:20:VAL:O	41:BS:23:LEU:HB2	2.16	0.45
43:DU:68:HIS:ND1	43:DU:70:SER:HB3	2.31	0.45
1:AA:934:C:H5	1:AA:1344:C:H2'	1.81	0.45
23:DA:519:U:H2'	23:DA:520:G:H8	1.82	0.45
32:DJ:105:LEU:O	32:DJ:106:LYS:C	2.55	0.45
23:DA:1461:G:O2'	23:DA:1462:C:H5'	2.16	0.45
42:BT:66:LEU:HD23	42:BT:67:GLY:N	2.30	0.45
23:BA:238:C:O2'	23:BA:608:A:H1'	2.16	0.45
25:BC:124:PRO:HG2	25:BC:129:ASN:ND2	2.31	0.45
2:CB:73:THR:HA	2:CB:94:ASN:O	2.16	0.45
23:BA:1427:A:H4'	23:BA:1428:C:O5'	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BM:51:ARG:O	35:BM:55:VAL:HG13	2.16	0.45
23:DA:2015:A:N3	50:D2:2:ALA:N	2.64	0.45
50:B2:4:HIS:HB2	50:B2:5:PRO:HD3	1.99	0.45
23:BA:1971:A:N3	25:BC:239:ARG:O	2.50	0.45
42:BT:25:LYS:HE3	42:BT:82:GLN:OE1	2.16	0.45
23:BA:1021:A:H2'	23:BA:1023:U:H5'	1.98	0.45
51:B3:25:LYS:HD3	53:B5:34:TRP:CZ3	2.51	0.45
23:DA:1021:A:H2'	23:DA:1023:U:H5'	1.97	0.45
34:DL:85:LEU:HA	34:DL:88:LEU:HB2	1.98	0.45
5:AE:50:GLU:HG3	5:AE:52:PRO:HD2	1.98	0.45
44:BV:71:VAL:HG11	44:BV:74:VAL:CG2	2.46	0.45
53:D5:54:GLU:O	53:D5:58:ILE:HG12	2.16	0.45
23:DA:2033:A:O2'	23:DA:2034:U:P	2.74	0.45
48:DZ:28:LEU:HA	48:DZ:33:GLN:OE1	2.16	0.45
34:DL:140:ALA:O	34:DL:141:ALA:CB	2.63	0.45
43:DU:9:LYS:O	43:DU:27:VAL:HG21	2.17	0.45
33:DK:71:ARG:HH21	33:DK:77:ILE:HG21	1.80	0.45
23:BA:2516:G:O6	23:BA:2517:C:N4	2.49	0.45
23:BA:2542:A:C8	23:BA:2544:G:O6	2.70	0.45
34:DL:12:ALA:HB1	34:DL:16:ARG:HB2	1.98	0.45
23:DA:861:A:H2'	23:DA:862:G:O4'	2.17	0.45
23:DA:2261:C:H3'	45:DW:16:SER:CB	2.46	0.45
28:BF:125:PHE:C	28:BF:127:GLY:H	2.20	0.45
40:DR:78:LYS:HG3	40:DR:79:VAL:HG23	1.98	0.45
13:CM:14:ARG:HG2	13:CM:44:ARG:NH1	2.32	0.45
22:CV:6170:G:H2'	22:CV:6171:U:C6	2.51	0.45
22:AV:6170:G:H2'	22:AV:6171:U:C6	2.52	0.45
23:DA:2271:G:H2'	23:DA:2272:U:H6	1.81	0.45
9:CI:17:VAL:HG21	9:CI:80:GLY:C	2.37	0.45
1:AA:947:G:H2'	1:AA:948:C:C6	2.51	0.45
38:BP:30:VAL:HG12	38:BP:86:ILE:HG12	1.99	0.45
23:BA:226:G:N2	23:BA:227:A:C2	2.85	0.45
23:DA:2563:U:H4'	33:DK:28:SER:HA	1.99	0.45
1:CA:191(G):G:C6	1:CA:192:U:C4	3.04	0.45
1:AA:191(G):G:C6	1:AA:192:U:C4	3.04	0.45
8:CH:35:ILE:O	8:CH:39:LEU:HB2	2.16	0.45
23:DA:2259:G:C2	23:DA:2282:G:N1	2.84	0.45
23:DA:2094:G:P	30:DH:22:LYS:HD2	2.57	0.45
25:BC:174:ILE:HD12	25:BC:174:ILE:N	2.31	0.45
30:BH:31:LEU:HD13	30:BH:37:VAL:HA	1.98	0.45
28:DF:39:ILE:HG12	28:DF:157:ILE:HG22	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BA:1178:C:H2'	23:BA:1179:C:C6	2.51	0.45
23:BA:1973:G:H2'	23:BA:1974:C:C6	2.51	0.45
2:CB:138:LEU:O	2:CB:141:GLU:HB2	2.16	0.45
41:BS:20:VAL:HG11	41:BS:44:ALA:HA	1.98	0.45
23:DA:2734:A:C8	23:DA:2735:G:C8	3.05	0.45
1:CA:902:G:H2'	1:CA:903:G:H8	1.82	0.45
23:DA:2531:A:H2	23:DA:2658:C:O2	1.99	0.45
1:AA:495:A:H4'	1:AA:496:A:OP1	2.16	0.45
38:DP:63:VAL:O	38:DP:73:GLU:HA	2.16	0.45
1:CA:384:G:H2'	1:CA:385:C:C6	2.51	0.45
3:CC:8:ILE:CD1	3:CC:16:ARG:HH21	2.28	0.45
23:DA:2764:A:N7	23:DA:2766:G:C6	2.84	0.45
32:DJ:101:TYR:N	32:DJ:101:TYR:CD1	2.85	0.45
1:CA:1031:G:H8	1:CA:1031:G:O5'	2.00	0.45
7:AG:45:ASP:O	7:AG:49:ILE:HG12	2.16	0.45
23:DA:1733:G:H8	23:DA:1733:G:O5'	2.00	0.45
26:BD:192:ASN:HD22	26:BD:192:ASN:N	2.15	0.45
23:DA:1208:C:C4	23:DA:1209:G:N7	2.85	0.45
23:DA:1954:G:N2	23:DA:1956:U:H3	2.14	0.45
34:BL:38:GLN:HG3	34:BL:41:ARG:HG2	1.99	0.45
39:DQ:92:ARG:HG2	40:DR:11:GLN:CD	2.37	0.45
29:BG:149:ARG:HD2	29:BG:164:TYR:HE1	1.82	0.45
1:AA:675:A:H2'	1:AA:676:A:H8	1.81	0.45
25:DC:27:THR:O	25:DC:27:THR:CG2	2.63	0.45
25:DC:32:SER:O	25:DC:33:LEU:O	2.35	0.45
5:AE:43:LEU:CD1	5:AE:132:ALA:HB1	2.41	0.45
7:CG:15:ASP:HB3	7:CG:19:GLY:N	2.28	0.45
1:AA:392:G:C2	1:AA:393:A:C4	3.05	0.45
23:BA:390:A:C5	34:BL:71:VAL:HG21	2.52	0.45
23:BA:1858:G:H1'	23:BA:1884:A:H61	1.82	0.45
5:AE:102:ALA:HB2	5:AE:120:THR:OG1	2.16	0.45
20:CT:72:LEU:HD23	20:CT:72:LEU:C	2.37	0.45
1:CA:973:G:H3'	1:CA:974:A:H5''	1.97	0.45
23:BA:186:G:H2'	23:BA:187:G:H8	1.82	0.45
23:DA:639:U:H2'	23:DA:640:C:H6	1.81	0.45
27:BE:83:PHE:O	27:BE:86:GLY:N	2.45	0.45
23:BA:36:G:N1	23:BA:445:C:C4	2.84	0.45
19:AS:10:PHE:O	19:AS:11:VAL:HB	2.16	0.45
23:DA:1587:A:H2'	23:DA:1588:C:H6	1.81	0.45
23:DA:813:U:H2'	23:DA:814:C:C6	2.52	0.45
23:DA:2306:C:H4'	28:DF:136:ARG:HH22	1.82	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
27:DE:124:LEU:HD12	27:DE:125:LEU:N	2.31	0.45
1:CA:687:A:H2'	1:CA:701:C:H41	1.82	0.45
1:CA:624:C:H4'	16:CP:10:GLY:HA2	1.98	0.45
23:BA:919:G:C5'	24:BB:81:G:H1'	2.46	0.45
14:CN:37:PHE:HE1	14:CN:53:LEU:HD22	1.82	0.45
6:AF:17:SER:O	6:AF:21:LEU:HD23	2.16	0.45
27:DE:29:ASN:N	27:DE:112:MET:HE1	2.32	0.45
19:CS:33:THR:HG22	19:CS:51:VAL:HA	1.98	0.45
1:AA:515:G:C2	1:AA:537:G:C2	3.05	0.45
23:DA:1871:A:H2'	23:DA:1872:A:C8	2.51	0.45
23:DA:1506:C:H2'	23:DA:1508:A:C8	2.50	0.45
23:DA:1229:G:H2'	23:DA:1230:C:C6	2.51	0.45
1:AA:191(F):U:O2	20:AT:105:SER:HB2	2.17	0.45
23:DA:1599:C:OP2	42:DT:36:LYS:HD3	2.17	0.45
1:AA:565:U:C6	1:AA:566:G:C8	3.05	0.45
9:CI:48:GLU:N	9:CI:49:PRO:CD	2.79	0.45
41:BS:70:TYR:O	41:BS:107:LEU:HA	2.17	0.45
1:AA:775:G:H2'	1:AA:776:G:O4'	2.16	0.45
1:CA:1106:G:H5''	3:CC:172:ARG:HG2	1.98	0.45
2:CB:32:ILE:HG12	2:CB:40:HIS:HD2	1.81	0.45
23:BA:868:U:C4	23:BA:869:G:N7	2.85	0.45
4:AD:201:GLN:O	4:AD:205:GLU:HG3	2.17	0.45
1:CA:1015:A:O5'	1:CA:1015:A:H8	1.99	0.45
1:CA:315:A:H5''	1:CA:317:G:OP2	2.17	0.45
12:CL:110:LYS:O	12:CL:111:ASP:HB2	2.16	0.45
1:AA:632:A:H2'	1:AA:633:G:O4'	2.16	0.45
1:AA:544:G:H2'	1:AA:545:C:C6	2.51	0.45
23:DA:1826:G:H2'	23:DA:1827:C:H6	1.81	0.45
43:BU:75:ILE:HG13	43:BU:80:GLY:H	1.82	0.45
37:DO:72:ALA:O	37:DO:76:LYS:HG3	2.17	0.45
3:CC:131:ARG:HH21	5:CE:50:GLU:HG2	1.81	0.45
38:DP:27:THR:HG22	38:DP:90:GLN:HB3	1.99	0.45
1:AA:942:G:H2'	1:AA:943:U:C6	2.52	0.45
34:DL:57:THR:HG23	34:DL:59:LEU:HB3	1.97	0.45
27:DE:53:THR:C	27:DE:55:GLY:N	2.70	0.45
40:BR:77:ALA:C	40:BR:79:VAL:H	2.19	0.45
23:BA:558:G:P	32:BJ:134:PRO:HD2	2.57	0.45
4:CD:8:VAL:C	4:CD:10:ARG:N	2.70	0.45
23:DA:479:A:HO2'	23:DA:481:G:H8	1.61	0.45
38:DP:50:ILE:HA	38:DP:99:LEU:CD1	2.46	0.45
23:DA:2807:G:C6	23:DA:2893:G:O6	2.69	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BA:777:A:O2'	23:BA:778:G:H5'	2.16	0.45
23:DA:2815:C:O2'	50:D2:42:PRO:HB2	2.16	0.45
38:BP:20:PRO:CD	38:BP:86:ILE:HG23	2.46	0.45
25:BC:70:TRP:CD1	25:BC:70:TRP:C	2.90	0.45
34:DL:27:HIS:CE1	40:DR:83:ARG:HH12	2.34	0.45
23:BA:1567:A:C8	25:BC:84:TYR:CE2	3.05	0.45
1:AA:501:C:OP1	12:AL:116:ARG:NH2	2.49	0.45
33:DK:31:LYS:HB3	33:DK:32:TYR:CE1	2.51	0.45
39:DQ:104:GLN:HB3	40:DR:44:LYS:NZ	2.32	0.45
1:AA:192:U:H2'	1:AA:193:C:H6	1.81	0.45
1:CA:642:A:HO2'	8:CH:31:PHE:HE1	1.65	0.45
23:DA:499:U:C4'	43:DU:47:LYS:HZ1	2.30	0.45
23:DA:2334:G:H4'	23:DA:2335:A:OP2	2.17	0.45
19:CS:53:ASN:C	19:CS:53:ASN:HD22	2.19	0.45
21:CU:14:TRP:CE3	21:CU:15:ARG:HG2	2.52	0.45
23:DA:116:C:O2'	23:DA:117:G:H5'	2.17	0.45
23:BA:2817:G:H21	23:BA:2836:U:H1'	1.82	0.45
23:BA:2722:G:H2'	23:BA:2723:C:C6	2.51	0.45
29:BG:38:SER:HB2	29:BG:41:MET:CG	2.46	0.45
23:DA:1272:A:O2'	23:DA:1273:U:H5'	2.17	0.45
23:DA:1599:C:H2'	23:DA:1600:C:C6	2.52	0.45
23:BA:2655:G:N2	23:BA:2664:G:C5	2.85	0.45
23:BA:553:U:O2'	23:BA:554:U:H5'	2.16	0.45
7:AG:86:GLN:HB2	7:AG:148:ASN:HD22	1.82	0.45
11:CK:48:ILE:HD11	11:CK:64:ALA:HA	1.99	0.45
26:BD:21:VAL:HG12	26:BD:23:VAL:HG13	1.99	0.45
49:B1:40:ILE:O	49:B1:47:VAL:HA	2.16	0.45
6:AF:50:TYR:CE1	18:AR:77:GLY:HA2	2.52	0.45
1:CA:102:G:H2'	1:CA:103:C:C6	2.51	0.45
23:BA:2436:G:C5	23:BA:2437:U:C5	3.04	0.45
1:CA:308:C:O2	1:CA:308:C:H2'	2.17	0.45
26:DD:203:LYS:HD2	26:DD:203:LYS:O	2.17	0.45
11:CK:108:ILE:O	18:CR:87:ARG:HA	2.16	0.45
26:DD:37:ARG:HD3	26:DD:42:ASP:OD1	2.17	0.45
1:AA:1454:G:H2'	1:AA:1455:G:H8	1.82	0.45
50:B2:4:HIS:HB2	50:B2:5:PRO:CD	2.46	0.45
23:DA:996:A:H4'	39:DQ:92:ARG:CZ	2.44	0.45
23:DA:1971:A:N3	25:DC:239:ARG:O	2.49	0.45
23:DA:1022:G:C6	23:DA:1140:C:C4	3.04	0.45
13:CM:102:ARG:HB3	13:CM:105:THR:OG1	2.17	0.45
28:DF:41:GLN:HG2	28:DF:155:MET:CB	2.37	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:BC:35:LYS:HD2	25:BC:35:LYS:HA	1.62	0.45
23:BA:2025:C:C2	23:BA:2026:C:C5	3.05	0.45
23:DA:2210:G:C3'	23:DA:2210:G:N3	2.77	0.45
23:BA:1006:C:H1'	32:BJ:129:MET:HG2	1.99	0.45
26:BD:201:THR:CG2	26:BD:202:LYS:H	2.22	0.45
41:BS:1:MET:HG3	41:BS:2:GLU:N	2.32	0.45
30:BH:9:LEU:HB2	30:BH:12:LEU:HB2	1.98	0.45
36:BN:10:LEU:HB2	36:BN:17:ARG:NH2	2.31	0.45
23:BA:1188:U:H4'	40:BR:79:VAL:HG13	1.98	0.45
23:BA:557:U:H2'	23:BA:558:G:C8	2.51	0.45
29:DG:46:GLU:HG3	29:DG:51:ARG:CD	2.47	0.45
12:AL:44:PRO:CD	12:AL:50:ALA:H	2.29	0.45
1:AA:1296:C:H5''	1:AA:1297:C:OP2	2.16	0.45
23:DA:644:A:C2	23:DA:646:A:C4	3.05	0.45
23:DA:1188:U:H4'	40:DR:79:VAL:HG13	1.99	0.45
23:BA:1348:G:C2'	23:BA:1349:A:H5''	2.45	0.45
23:DA:1804:C:O5'	23:DA:1804:C:H6	1.99	0.45
23:DA:1694:C:C5'	23:DA:1694:C:C6	2.96	0.45
25:BC:11:PRO:C	25:BC:13:ARG:N	2.70	0.45
25:DC:211:ARG:O	25:DC:215:LEU:HG	2.16	0.45
1:AA:683:G:C5	1:AA:684:A:N7	2.85	0.45
23:BA:2807:G:C6	23:BA:2893:G:O6	2.69	0.45
1:CA:505:G:C6	1:CA:535:A:C2	3.04	0.45
23:BA:2346:A:H5''	23:BA:2383:G:C1'	2.46	0.45
33:DK:22:ILE:HD13	33:DK:22:ILE:HA	1.77	0.45
29:BG:13:LYS:HE2	29:BG:14:GLY:N	2.32	0.45
1:CA:1075:C:H5''	2:CB:179:LYS:HZ3	1.79	0.45
1:CA:515:G:C2	1:CA:537:G:C2	3.04	0.45
23:BA:1952:A:C4	33:BK:22:ILE:HD12	2.50	0.45
1:CA:554:C:H2'	1:CA:555:C:H6	1.82	0.45
1:CA:192:U:O2'	1:CA:193:C:H5'	2.17	0.45
14:AN:37:PHE:HE1	14:AN:53:LEU:HD22	1.81	0.45
1:AA:1410:G:H2'	1:AA:1411:C:H6	1.81	0.45
23:BA:356:G:H2'	23:BA:357:A:H8	1.81	0.45
23:DA:2817:G:H21	23:DA:2836:U:H1'	1.82	0.45
6:CF:50:TYR:CE1	18:CR:77:GLY:HA2	2.52	0.45
1:CA:1250:A:H5'	9:CI:67:GLY:HA2	1.98	0.45
1:CA:342:C:N3	1:CA:348:G:C2	2.85	0.45
47:DY:31:GLU:O	47:DY:35:LEU:HB2	2.17	0.45
2:AB:32:ILE:HG12	2:AB:40:HIS:HD2	1.81	0.45
32:DJ:26:THR:HG22	32:DJ:27:TYR:N	2.31	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BA:1489:U:HO2'	23:BA:1490:A:H8	1.65	0.45
23:BA:1208:C:C4	23:BA:1209:G:N7	2.85	0.45
1:AA:1318:A:O2'	19:AS:37:ARG:HB2	2.16	0.45
44:DV:152:ALA:N	44:DV:169:GLU:O	2.48	0.45
2:CB:153:ARG:NH1	2:CB:153:ARG:HB2	2.31	0.45
16:AP:43:LYS:HA	16:AP:48:TRP:HB3	1.99	0.45
38:DP:115:ARG:H	38:DP:115:ARG:HG2	1.53	0.45
23:DA:1850:G:C6	23:DA:1851:U:C4	3.05	0.45
32:DJ:93:LYS:HB3	32:DJ:110:LEU:HB2	1.99	0.45
9:AI:39:GLY:O	9:AI:40:LEU:HD23	2.17	0.45
23:BA:2552:U:H2'	23:BA:2554:U:OP2	2.16	0.45
40:DR:38:LEU:C	40:DR:39:LEU:HD22	2.36	0.45
23:DA:2394:C:H2'	23:DA:2395:C:H6	1.81	0.45
1:AA:735:C:C2	1:AA:736:C:C5	3.05	0.45
45:BW:32:ARG:C	45:BW:35:ASN:HD21	2.19	0.45
28:BF:106:LEU:HB3	28:BF:107:LEU:HD23	1.99	0.45
5:AE:152:ARG:HD3	8:AH:42:GLU:O	2.16	0.45
27:BE:53:THR:C	27:BE:55:GLY:N	2.70	0.45
1:CA:321:A:N7	1:CA:328:C:C6	2.84	0.45
26:DD:6:GLY:HA2	26:DD:51:PHE:CZ	2.51	0.45
43:BU:27:VAL:O	43:BU:27:VAL:HG23	2.16	0.45
33:DK:77:ILE:HD13	33:DK:78:ARG:N	2.31	0.45
26:BD:6:GLY:HA2	26:BD:51:PHE:CZ	2.51	0.45
23:DA:661:C:H2'	23:DA:662:G:H8	1.82	0.45
23:DA:643:A:C2	23:DA:644:A:C4	3.05	0.45
3:AC:20:SER:HB2	3:AC:40:ARG:NH1	2.31	0.45
51:D3:34:LEU:HD13	51:D3:34:LEU:N	2.32	0.45
1:CA:406:G:N2	1:CA:437:U:C2	2.85	0.45
11:AK:33:THR:HA	11:AK:40:ILE:HG12	1.98	0.45
32:BJ:116:THR:HG23	32:BJ:117:HIS:N	2.32	0.45
9:AI:5:TYR:HA	9:AI:17:VAL:O	2.17	0.45
2:AB:168:THR:HG1	2:AB:192:SER:HA	1.81	0.45
1:AA:1298:C:H41	7:AG:114:ARG:HD3	1.81	0.45
42:DT:64:LYS:HG2	42:DT:65:ARG:N	2.30	0.45
6:CF:61:LEU:HD23	6:CF:63:TYR:OH	2.17	0.45
8:CH:97:VAL:HG13	8:CH:98:LYS:H	1.81	0.45
43:DU:75:ILE:HG12	43:DU:76:CYS:N	2.32	0.45
23:BA:2850:A:H8	23:BA:2850:A:H5'	1.79	0.45
23:BA:301:G:C6	23:BA:317:G:C6	3.05	0.45
1:AA:1501:C:C2	1:AA:1504:G:O6	2.70	0.45
1:AA:1502:A:C8	1:AA:1505:G:N2	2.85	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:818:G:C3'	1:CA:819:A:H5''	2.47	0.45
1:AA:1281:U:H5'	1:AA:1282:C:H5	1.82	0.45
37:DO:38:GLN:HB3	37:DO:47:THR:HG23	1.98	0.45
46:DX:21:ARG:HD3	46:DX:21:ARG:HA	1.86	0.45
23:DA:380:U:H4'	46:DX:21:ARG:O	2.17	0.45
23:DA:1615:C:C6	23:DA:1617:C:C5	3.04	0.45
26:DD:78:LEU:C	26:DD:79:ARG:HD2	2.36	0.45
1:AA:645:C:H2'	1:AA:646:U:O4'	2.16	0.45
23:DA:1326:U:O2'	23:DA:1327:C:H5'	2.17	0.45
23:BA:2580:U:H4'	26:BD:130:GLY:HA2	1.98	0.45
1:AA:1396:A:C2	5:AE:19:MET:HG3	2.52	0.45
26:DD:24:THR:HB	26:DD:186:GLY:O	2.17	0.45
26:BD:170:LEU:HB3	26:BD:184:VAL:HG12	1.99	0.45
1:CA:1468:A:H2'	1:CA:1469:G:O4'	2.17	0.45
1:CA:284:G:H2'	1:CA:285:G:H8	1.80	0.45
1:AA:862:C:O2'	1:AA:863:U:H5'	2.17	0.45
23:DA:2515:C:H1'	23:DA:2570:G:N2	2.32	0.45
1:CA:843:U:H5'	1:CA:848:C:C6	2.52	0.45
42:BT:62:LYS:O	42:BT:73:ARG:HB2	2.16	0.45
23:BA:1682:G:H2'	23:BA:1683:C:C6	2.51	0.45
2:AB:153:ARG:NH1	2:AB:153:ARG:HB2	2.31	0.45
1:CA:695:A:H2'	1:CA:696:A:C8	2.52	0.45
23:BA:979:G:H3'	23:BA:980:A:H5''	1.99	0.45
23:BA:244:A:C2	23:BA:255:A:C4	3.05	0.45
23:BA:994:C:OP1	39:BQ:53:ARG:NH2	2.49	0.45
34:DL:128:HIS:HB3	34:DL:147:LEU:HD23	1.97	0.45
47:BY:33:MET:O	47:BY:37:PHE:HB2	2.17	0.45
23:BA:1022:G:C5	23:BA:1140:C:N4	2.85	0.45
23:BA:142:G:H1'	42:BT:37:THR:HG21	1.98	0.45
27:BE:63:LYS:NZ	27:BE:67:GLN:HE21	2.15	0.45
23:BA:2712:U:O2'	23:BA:712(B):A:P	2.75	0.45
5:CE:50:GLU:HG3	5:CE:52:PRO:HD2	1.98	0.45
23:BA:1578:U:C2'	23:BA:1579:A:H5''	2.46	0.45
5:CE:43:LEU:HD12	5:CE:109:ILE:HD11	1.99	0.45
28:BF:88:ILE:HG13	28:BF:89:GLY:N	2.30	0.45
17:CQ:69:LYS:O	17:CQ:70:ARG:HD2	2.16	0.45
27:DE:160:ASN:ND2	27:DE:162:LEU:H	2.15	0.45
11:AK:20:TYR:C	11:AK:21:ILE:HD12	2.37	0.45
4:CD:30:LYS:HD3	4:CD:35:ARG:NH2	2.32	0.45
3:CC:6:HIS:CD2	3:CC:7:PRO:HD2	2.43	0.45
23:BA:66:C:H2'	23:BA:67:U:H6	1.81	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
27:DE:89:VAL:HG12	27:DE:90:PHE:CD2	2.52	0.45
13:AM:30:ALA:O	13:AM:34:LEU:HG	2.17	0.45
9:AI:17:VAL:HG22	9:AI:63:ILE:HG13	1.99	0.45
23:BA:598:G:H2'	23:BA:599:G:O4'	2.16	0.45
32:BJ:127:LYS:HB2	32:BJ:140:PHE:HE1	1.78	0.45
23:BA:1291:C:H2'	23:BA:1292:U:H6	1.79	0.45
25:BC:5:LYS:H	25:BC:5:LYS:HD2	1.80	0.45
43:BU:90:LEU:HD23	43:BU:90:LEU:N	2.31	0.45
25:DC:70:TRP:C	25:DC:70:TRP:CD1	2.90	0.45
10:AJ:6:ILE:HG22	10:AJ:98:ILE:CG2	2.45	0.45
29:DG:13:LYS:HE2	29:DG:14:GLY:N	2.32	0.45
2:CB:112:VAL:O	2:CB:115:LEU:HB3	2.17	0.45
1:AA:262:A:C6	1:AA:263:A:C6	3.05	0.45
34:BL:105:LEU:HD12	34:BL:105:LEU:H	1.82	0.45
44:DV:8:TYR:HB2	44:DV:38:TYR:CE2	2.52	0.45
22:AV:6167:G:H2'	22:AV:6168:G:C8	2.52	0.45
1:CA:740:U:O3'	15:CO:39:LEU:HD23	2.16	0.45
19:AS:53:ASN:HD22	19:AS:53:ASN:C	2.20	0.45
45:BW:50:ASN:ND2	45:BW:83:PRO:HD3	2.31	0.45
23:BA:454:A:H4'	23:BA:455:C:OP2	2.17	0.45
15:CO:62:GLN:O	15:CO:66:LEU:HD13	2.17	0.45
23:DA:2748:A:C2	23:DA:2757:A:C4	3.04	0.45
1:CA:246:A:C2	1:CA:282:A:C5	3.05	0.45
29:BG:105:LEU:N	29:BG:105:LEU:HD13	2.31	0.45
7:CG:86:GLN:HB2	7:CG:148:ASN:HD22	1.81	0.45
4:CD:103:ASN:OD1	4:CD:114:ARG:NH2	2.50	0.45
23:DA:311:A:C8	23:DA:332:A:C5	3.05	0.45
34:BL:50:ARG:HB3	53:B5:60:LEU:HD21	1.98	0.45
34:DL:39:LYS:HA	34:DL:39:LYS:HD3	1.71	0.45
1:CA:1130:A:N6	1:CA:1144:G:H21	2.15	0.45
32:DJ:42:GLU:O	32:DJ:42:GLU:HG3	2.17	0.45
23:DA:637:A:OP2	34:DL:115:LEU:HB2	2.16	0.45
35:BM:85:LYS:HG3	35:BM:86:GLY:N	2.32	0.45
23:DA:846:C:C2	23:DA:847:U:C5	3.05	0.45
26:BD:91:VAL:HB	26:BD:95:ILE:CD1	2.40	0.45
53:D5:14:VAL:HG22	53:D5:24:ALA:HB2	1.99	0.45
1:CA:81:G:C5	1:CA:82:U:C4	3.05	0.45
23:BA:1006:C:C2	23:BA:1138:G:N2	2.85	0.45
23:BA:1658:C:OP1	26:BD:132:HIS:O	2.34	0.45
23:BA:2543:G:H2'	23:BA:2544:G:O4'	2.17	0.45
23:BA:2287:A:C4	23:BA:2289:G:C8	3.04	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:DL:16:ARG:NH2	34:DL:18:ARG:H	2.15	0.45
35:BM:101:ARG:HG3	35:BM:102:VAL:N	2.32	0.45
32:DJ:116:THR:O	32:DJ:118:PRO:HD3	2.17	0.45
5:AE:90:VAL:O	5:AE:120:THR:HA	2.17	0.45
44:BV:22:GLY:O	44:BV:41:LEU:HG	2.17	0.45
23:BA:1478:G:C2	23:BA:1479:G:C8	3.04	0.45
23:BA:1390:U:O2'	23:BA:1391:U:H5'	2.16	0.45
23:DA:2346:A:C2	23:DA:2383:G:C2	3.05	0.45
23:BA:2850:A:H5''	23:BA:2868:A:C2	2.51	0.45
32:BJ:62:ARG:NE	32:BJ:64:ASP:HB2	2.32	0.45
1:CA:324:G:N2	1:CA:327:A:C8	2.85	0.45
25:DC:108:PRO:HB3	25:DC:143:HIS:CE1	2.52	0.45
27:BE:124:LEU:HB3	27:BE:193:VAL:HG22	1.99	0.45
1:AA:740:U:O3'	15:AO:39:LEU:HD23	2.16	0.45
27:DE:192:LEU:C	27:DE:192:LEU:HD23	2.37	0.45
1:CA:818:G:H3'	1:CA:819:A:H5''	1.98	0.45
1:CA:99:C:C6	1:CA:99:C:C3'	2.99	0.45
23:BA:380:U:H4'	46:BX:21:ARG:O	2.17	0.45
1:CA:1305:G:C8	1:CA:1305:G:OP2	2.70	0.45
23:BA:2196:C:O2'	23:BA:2197:U:H5'	2.17	0.45
1:CA:964:A:H5'	1:CA:1199:U:OP1	2.16	0.45
23:DA:155:C:H2'	23:DA:161:U:H5'	1.99	0.45
23:BA:2259:G:C2	23:BA:2282:G:N1	2.85	0.45
23:BA:2886:G:H2'	23:BA:2887:U:H6	1.82	0.45
23:DA:1600:C:O2'	23:DA:1601:G:H5'	2.17	0.45
35:DM:51:ARG:O	35:DM:55:VAL:HG13	2.18	0.45
51:D3:41:PRO:HG3	51:D3:49:HIS:HE1	1.82	0.45
1:AA:284:G:H2'	1:AA:285:G:H8	1.82	0.45
25:DC:124:PRO:HG2	25:DC:129:ASN:ND2	2.32	0.45
23:DA:1427:A:H4'	23:DA:1428:C:O5'	2.17	0.45
1:CA:179:A:H2'	1:CA:180:U:H6	1.81	0.45
1:AA:1057:G:H4'	3:AC:197:GLY:N	2.32	0.45
23:BA:2315:G:H2'	23:BA:2316:C:C6	2.52	0.45
25:DC:218:ARG:HH11	25:DC:218:ARG:HG3	1.82	0.45
27:BE:54:ARG:HA	27:BE:87:GLY:HA3	1.98	0.45
42:DT:55:ASN:HB2	42:DT:80:ILE:HG23	1.99	0.44
32:BJ:156:GLN:O	32:BJ:157:ARG:HB2	2.17	0.44
12:CL:44:PRO:HG3	12:CL:52:ARG:CD	2.46	0.44
35:DM:22:LYS:CE	35:DM:22:LYS:HA	2.35	0.44
43:DU:81:LYS:NZ	43:DU:99:CYS:SG	2.90	0.44
23:DA:2250:G:C4	35:DM:82:ARG:HG3	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
38:DP:88:ILE:HG13	38:DP:89:VAL:N	2.31	0.44
28:BF:58:GLN:O	28:BF:62:LEU:HD13	2.17	0.44
5:CE:33:VAL:CG1	5:CE:109:ILE:HD13	2.47	0.44
48:DZ:26:LEU:HB2	48:DZ:28:LEU:CD1	2.47	0.44
23:DA:1578:U:C2'	23:DA:1579:A:H5''	2.47	0.44
30:BH:82:ARG:HD2	30:BH:89:TYR:CD2	2.52	0.44
12:CL:29:ALA:HA	12:CL:30:PRO:HD3	1.78	0.44
23:BA:2037:G:C6	23:BA:2038:G:C6	3.05	0.44
23:DA:2287:A:C4	23:DA:2289:G:C8	3.04	0.44
32:BJ:57:LEU:CD1	32:BJ:142:ARG:HB2	2.48	0.44
23:DA:662:G:OP1	34:DL:18:ARG:HD2	2.17	0.44
23:DA:1448:G:N2	23:DA:149(B):A:N6	2.66	0.44
37:BO:90:GLY:O	37:BO:92:TYR:CD1	2.69	0.44
28:DF:125:PHE:C	28:DF:127:GLY:H	2.19	0.44
23:DA:644:A:O2'	23:DA:645:C:H5''	2.17	0.44
7:AG:15:ASP:CB	7:AG:20:ASP:H	2.29	0.44
13:CM:30:ALA:O	13:CM:34:LEU:HG	2.17	0.44
25:DC:11:PRO:C	25:DC:13:ARG:N	2.70	0.44
12:CL:84:ILE:HD12	12:CL:84:ILE:N	2.32	0.44
9:CI:5:TYR:HA	9:CI:17:VAL:O	2.16	0.44
25:BC:211:ARG:O	25:BC:215:LEU:HG	2.17	0.44
29:DG:105:LEU:HD13	29:DG:105:LEU:N	2.31	0.44
42:BT:64:LYS:HG2	42:BT:65:ARG:N	2.31	0.44
44:DV:58:VAL:HG11	44:DV:66:SER:HB2	1.98	0.44
23:DA:2666:C:H3'	23:DA:2667:C:C6	2.52	0.44
1:AA:506:G:C6	1:AA:507:C:C4	3.04	0.44
39:BQ:104:GLN:HB3	40:BR:44:LYS:NZ	2.32	0.44
23:BA:102:G:H5''	23:BA:102:G:H8	1.82	0.44
13:CM:27:LYS:HE2	13:CM:31:LYS:CE	2.46	0.44
23:DA:813:U:O2'	23:DA:1225:G:H1'	2.17	0.44
1:CA:262:A:H5'	20:CT:74:LYS:HG3	1.98	0.44
33:BK:86:ILE:HD12	33:BK:86:ILE:N	2.31	0.44
1:AA:554:C:H2'	1:AA:555:C:H6	1.82	0.44
53:B5:11:LYS:HB2	53:B5:61:LEU:HD22	1.99	0.44
28:DF:16:ARG:HB3	28:DF:17:PRO:HD3	1.98	0.44
23:DA:404:C:H4'	23:DA:405:U:H5'	1.99	0.44
4:AD:68:TYR:CE2	4:AD:97:LEU:HB3	2.51	0.44
15:CO:65:ARG:O	15:CO:68:ARG:HB2	2.17	0.44
23:BA:225:A:O2'	23:BA:257:A:H4'	2.17	0.44
29:DG:122:THR:O	29:DG:134:SER:HB2	2.16	0.44
41:DS:70:TYR:O	41:DS:107:LEU:HA	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
29:BG:24:VAL:HG23	29:BG:37:VAL:HG21	1.98	0.44
1:CA:862:C:O2'	1:CA:863:U:H5'	2.17	0.44
45:BW:27:GLU:HB2	45:BW:69:PHE:HD1	1.81	0.44
1:AA:695:A:H2'	1:AA:696:A:C8	2.52	0.44
26:BD:37:ARG:HD3	26:BD:42:ASP:OD1	2.18	0.44
16:AP:55:ARG:O	16:AP:58:TYR:HB3	2.17	0.44
7:AG:107:ALA:HB2	7:AG:134:ALA:HB2	1.99	0.44
8:CH:13:ILE:O	8:CH:17:THR:HG23	2.16	0.44
8:AH:1:MET:N	8:AH:1:MET:HE2	2.32	0.44
4:CD:201:GLN:O	4:CD:205:GLU:HG3	2.17	0.44
23:BA:1900:A:N1	23:BA:1970:A:C6	2.85	0.44
5:CE:77:PRO:HG2	5:CE:78:HIS:ND1	2.32	0.44
13:AM:102:ARG:HB3	13:AM:105:THR:OG1	2.17	0.44
23:BA:2250:G:C4	35:BM:82:ARG:HG3	2.52	0.44
49:B1:64:LYS:HA	49:B1:64:LYS:HE3	2.00	0.44
49:D1:46:ASN:HB2	49:D1:64:LYS:CB	2.47	0.44
23:DA:2712:U:O2'	23:DA:712(B):A:P	2.75	0.44
35:DM:141:GLN:HE21	44:DV:72:ARG:HG2	1.82	0.44
23:DA:744:G:OP1	26:DD:132:HIS:HB2	2.18	0.44
23:DA:572:A:H5''	23:DA:573:G:OP2	2.16	0.44
23:BA:2028:U:O4	23:BA:2033:A:OP1	2.35	0.44
1:CA:1014:A:H5'	19:CS:14:HIS:NE2	2.32	0.44
37:BO:18:ILE:HA	37:BO:21:THR:OG1	2.17	0.44
23:DA:1652:A:H2'	23:DA:1653:G:O4'	2.18	0.44
27:BE:139:PHE:HB2	27:BE:166:ALA:CB	2.45	0.44
23:BA:1317:A:C6	23:BA:1318:C:C4	3.05	0.44
34:BL:30:THR:HG22	34:BL:31:ALA:H	1.80	0.44
44:DV:137:ILE:HD12	44:DV:137:ILE:N	2.33	0.44
1:AA:1511:G:C6	1:AA:1512:U:C4	3.05	0.44
1:CA:1187:G:H2'	1:CA:1188:A:H8	1.82	0.44
46:DX:90:ILE:O	46:DX:94:LEU:N	2.50	0.44
26:DD:118:LYS:NZ	36:DN:2:ARG:HH22	2.15	0.44
26:BD:4:ILE:CG2	26:BD:96:PHE:HE1	2.30	0.44
38:DP:98:LYS:HB3	38:DP:100:TYR:HE1	1.82	0.44
1:AA:9:G:OP2	5:AE:121:LYS:HG3	2.17	0.44
23:BA:2809:A:C6	23:BA:2892:A:C8	3.05	0.44
5:AE:92:LYS:HB3	5:AE:119:LEU:HB2	1.99	0.44
1:AA:1528:U:H5''	1:AA:1528:U:C6	2.50	0.44
23:BA:1603:A:H5'	23:BA:1603:A:H8	1.81	0.44
29:DG:20:ALA:HB1	29:DG:21:PRO:CD	2.47	0.44
24:BB:111:U:HO2'	24:BB:112:G:H8	1.62	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
44:BV:92:SER:HB2	44:BV:94:GLU:OE2	2.16	0.44
1:CA:37:U:OP2	12:CL:122:LYS:HE3	2.18	0.44
15:CO:53:HIS:CE1	23:DA:715:G:O6	2.71	0.44
1:AA:1305:G:OP2	1:AA:1305:G:C8	2.71	0.44
29:DG:22:GLY:C	29:DG:23:ARG:HD3	2.38	0.44
1:AA:909:A:H3'	1:AA:910:C:H6	1.82	0.44
19:CS:58:VAL:O	19:CS:58:VAL:HG23	2.18	0.44
40:BR:99:ILE:H	40:BR:99:ILE:HD13	1.81	0.44
23:DA:1797:C:O2'	25:DC:259:THR:HG23	2.17	0.44
23:BA:1600:C:O2'	23:BA:1601:G:H5'	2.17	0.44
1:CA:284:G:H2'	1:CA:285:G:C8	2.52	0.44
23:BA:816:C:O2'	23:BA:817:C:H5'	2.16	0.44
1:AA:838:G:N2	1:AA:849:C:C2	2.85	0.44
30:BH:7:GLU:OE1	30:BH:8:PRO:HD2	2.16	0.44
23:BA:1916:A:H2'	23:BA:1917:U:O4'	2.18	0.44
44:BV:150:LEU:HD23	44:BV:151:HIS:N	2.32	0.44
23:BA:570:G:H2'	23:BA:2030:A:C5	2.52	0.44
7:CG:74:GLU:HG2	7:CG:91:VAL:HG22	2.00	0.44
1:CA:1207:G:H2'	1:CA:1208:C:H6	1.82	0.44
34:DL:67:MET:HE3	34:DL:67:MET:HA	1.99	0.44
49:D1:40:ILE:O	49:D1:47:VAL:HA	2.17	0.44
1:AA:179:A:H2'	1:AA:180:U:H6	1.82	0.44
9:AI:19:LEU:CD2	9:AI:59:PHE:HB3	2.30	0.44
23:DA:783:A:H3'	23:DA:783:A:C8	2.52	0.44
1:CA:1371:G:O3'	9:CI:69:GLY:HA3	2.17	0.44
34:BL:57:THR:C	34:BL:59:LEU:H	2.21	0.44
41:DS:1:MET:HG3	41:DS:2:GLU:N	2.32	0.44
25:DC:30:GLU:HG3	25:DC:63:ARG:CZ	2.47	0.44
5:CE:152:ARG:HD3	8:CH:42:GLU:O	2.17	0.44
10:AJ:74:ILE:HG12	10:AJ:74:ILE:O	2.17	0.44
1:CA:199:G:C2'	1:CA:200:G:H5''	2.42	0.44
28:DF:25:TYR:CZ	28:DF:32:PRO:HD3	2.51	0.44
4:AD:30:LYS:HD3	4:AD:35:ARG:NH2	2.32	0.44
30:DH:128:LEU:HG	30:DH:142:VAL:CG2	2.48	0.44
32:BJ:116:THR:O	32:BJ:118:PRO:HD3	2.18	0.44
32:DJ:114:LEU:HA	32:DJ:118:PRO:HB3	1.99	0.44
23:DA:1478:G:C2	23:DA:1479:G:C8	3.05	0.44
1:CA:801:U:H2'	1:CA:802:A:H8	1.80	0.44
8:AH:64:LYS:HD2	8:AH:79:VAL:HG11	1.97	0.44
26:DD:4:ILE:CG2	26:DD:96:PHE:HE1	2.29	0.44
32:DJ:62:ARG:NE	32:DJ:64:ASP:HB2	2.32	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DA:226:G:N2	23:DA:227:A:C2	2.84	0.44
28:BF:115:ARG:CD	28:BF:115:ARG:H	2.31	0.44
23:DA:1378:A:H4'	23:DA:1379:A:OP1	2.17	0.44
5:AE:12:LEU:C	5:AE:13:ILE:HD12	2.38	0.44
23:BA:17:G:H2'	23:BA:18:C:C6	2.52	0.44
1:CA:390:C:O3'	16:CP:28:ARG:NH2	2.50	0.44
23:BA:404:C:H4'	23:BA:405:U:H5'	1.99	0.44
4:CD:57:ARG:HB3	4:CD:206:PHE:HB2	1.99	0.44
23:DA:753:C:H2'	23:DA:754:C:H6	1.82	0.44
53:D5:11:LYS:HB2	53:D5:61:LEU:HD22	2.00	0.44
28:BF:174:GLU:HG2	28:BF:180:PHE:HD1	1.80	0.44
23:BA:1487:G:H2'	23:BA:1488:G:H8	1.81	0.44
9:AI:52:ALA:C	9:AI:95:LYS:HZ1	2.21	0.44
23:DA:792:G:H5''	23:DA:793:A:H5'	1.98	0.44
1:CA:1010:G:H2'	1:CA:1011:G:C8	2.53	0.44
1:CA:579:G:H2'	1:CA:580:U:C6	2.51	0.44
1:CA:1442:G:H8	1:CA:1442:G:H3'	1.81	0.44
5:AE:25:ARG:HD2	5:AE:25:ARG:N	2.33	0.44
15:AO:33:THR:HA	15:AO:63:ARG:HH12	1.82	0.44
42:DT:24:GLY:O	42:DT:83:VAL:HG22	2.17	0.44
2:CB:162:ILE:O	2:CB:185:ILE:HG12	2.17	0.44
23:BA:142:G:H1'	42:BT:37:THR:CG2	2.47	0.44
43:BU:15:VAL:HG22	43:BU:72:VAL:HG12	1.99	0.44
28:DF:58:GLN:O	28:DF:62:LEU:HD13	2.17	0.44
26:DD:91:VAL:HB	26:DD:95:ILE:CD1	2.39	0.44
35:BM:141:GLN:HE21	44:BV:72:ARG:HG2	1.82	0.44
36:DN:54:LEU:HD22	36:DN:66:VAL:HG23	1.99	0.44
28:DF:106:LEU:HB3	28:DF:107:LEU:HD23	1.99	0.44
23:DA:2027:G:C2	23:DA:2028:U:H1'	2.53	0.44
3:AC:150:LYS:O	3:AC:200:ALA:HA	2.17	0.44
16:CP:4:ILE:N	16:CP:4:ILE:HD12	2.31	0.44
36:DN:9:LYS:HD3	36:DN:43:GLU:OE1	2.18	0.44
3:CC:21:ARG:O	3:CC:58:GLU:HA	2.16	0.44
28:BF:172:LEU:O	28:BF:176:LEU:HG	2.17	0.44
19:AS:29:ARG:O	19:AS:31:ILE:HG22	2.16	0.44
4:AD:8:VAL:C	4:AD:10:ARG:N	2.70	0.44
12:AL:44:PRO:HG3	12:AL:52:ARG:CD	2.46	0.44
30:DH:114:LEU:HA	30:DH:130:TYR:HD1	1.83	0.44
32:BJ:116:THR:OG1	32:BJ:117:HIS:N	2.51	0.44
23:DA:2598:A:H5''	25:DC:235:GLY:HA2	1.98	0.44
10:CJ:55:LYS:O	10:CJ:56:HIS:CG	2.70	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
30:BH:6:LEU:HD23	30:BH:36:ALA:HA	1.99	0.44
1:AA:623:C:C4	1:AA:624:C:C5	3.04	0.44
6:CF:63:TYR:O	6:CF:65:VAL:HG13	2.18	0.44
23:BA:2849:U:H4'	23:BA:2868:A:C2	2.52	0.44
18:AR:56:THR:CB	18:AR:58:LEU:HD13	2.48	0.44
11:AK:57:THR:HG22	11:AK:59:TYR:N	2.33	0.44
23:DA:828:U:O2	23:DA:828:U:H3'	2.18	0.44
38:DP:30:VAL:HG12	38:DP:86:ILE:HG12	1.99	0.44
1:AA:236:G:C6	1:AA:237:C:C4	3.05	0.44
44:BV:6:LYS:HG3	44:BV:8:TYR:CZ	2.53	0.44
45:DW:51:VAL:HG21	45:DW:80:HIS:HA	2.00	0.44
3:AC:186:PHE:CG	3:AC:187:ALA:N	2.86	0.44
23:BA:363(C):G:H2'	23:BA:363(D):G:H8	1.83	0.44
23:DA:1536:A:H5"	23:DA:1537:C:OP2	2.17	0.44
44:DV:10:ARG:HG2	44:DV:11:GLU:N	2.32	0.44
1:CA:833:U:C2	1:CA:834:C:C5	3.05	0.44
18:CR:51:LEU:HA	18:CR:52:PRO:HD3	1.87	0.44
23:BA:357:A:H2'	23:BA:358:U:C6	2.52	0.44
25:DC:260:ARG:O	25:DC:261:LYS:C	2.55	0.44
38:DP:62:THR:HG22	38:DP:75:ILE:HG13	1.98	0.44
23:BA:2820:A:O4'	36:BN:5:LYS:HG3	2.17	0.44
23:DA:2330:G:O2'	45:DW:41:ARG:HB2	2.18	0.44
44:BV:54:HIS:CG	44:BV:101:PRO:HG3	2.53	0.44
41:BS:22:ASP:HA	41:BS:25:ARG:HH12	1.83	0.44
25:DC:175:LEU:HD12	25:DC:185:VAL:HG21	1.98	0.44
1:AA:409:G:OP1	4:AD:24:GLU:N	2.50	0.44
23:DA:616:A:C4'	23:DA:617:G:OP1	2.66	0.44
4:CD:128:VAL:HA	4:CD:145:GLU:O	2.16	0.44
30:DH:7:GLU:OE1	30:DH:8:PRO:HD2	2.17	0.44
23:DA:1916:A:H2'	23:DA:1917:U:O4'	2.18	0.44
23:DA:2501:C:H6	23:DA:2501:C:H2'	1.56	0.44
1:AA:945:G:H2'	1:AA:945:G:N3	2.33	0.44
1:AA:308:C:H2'	1:AA:308:C:O2	2.17	0.44
1:AA:1031:G:O5'	1:AA:1031:G:H8	2.00	0.44
29:DG:12:PRO:O	29:DG:15:VAL:HG22	2.18	0.44
17:AQ:7:THR:HG22	17:AQ:58:GLU:HG2	1.99	0.44
9:AI:45:ALA:O	9:AI:48:GLU:HB2	2.17	0.44
34:DL:38:GLN:CG	34:DL:39:LYS:H	2.28	0.44
23:BA:71:A:C2	42:BT:31:HIS:CE1	3.04	0.44
37:DO:98:VAL:O	37:DO:101:LEU:HB2	2.18	0.44
35:DM:141:GLN:N	44:DV:53:ILE:HB	2.33	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:BN:54:LEU:HD22	36:BN:66:VAL:HG23	2.00	0.44
23:BA:572:A:H5''	23:BA:573:G:OP2	2.17	0.44
23:DA:1174:A:H3'	23:DA:1175:U:C5'	2.46	0.44
30:BH:83:ALA:CA	30:BH:89:TYR:HD1	2.29	0.44
36:DN:10:LEU:HD23	36:DN:21:TYR:OH	2.18	0.44
23:DA:2484:G:C2	23:DA:2485:G:C8	3.06	0.44
7:CG:15:ASP:CB	7:CG:20:ASP:H	2.30	0.44
28:DF:172:LEU:O	28:DF:176:LEU:HG	2.18	0.44
3:CC:20:SER:HB2	3:CC:40:ARG:NH1	2.31	0.44
23:BA:1654:A:OP2	36:BN:3:HIS:NE2	2.51	0.44
29:BG:123:PHE:HA	29:BG:133:VAL:HA	1.99	0.44
43:DU:90:LEU:N	43:DU:90:LEU:HD23	2.33	0.44
23:BA:1030:G:OP2	35:BM:128:LYS:HE3	2.17	0.44
12:AL:84:ILE:HD12	12:AL:84:ILE:N	2.33	0.44
25:BC:108:PRO:HB3	25:BC:143:HIS:CE1	2.52	0.44
1:CA:236:G:C6	1:CA:237:C:C4	3.05	0.44
27:DE:124:LEU:HB3	27:DE:193:VAL:HG22	1.98	0.44
6:CF:62:TRP:CB	18:CR:35:ARG:HH12	2.30	0.44
1:CA:624:C:H4'	16:CP:11:SER:N	2.33	0.44
1:AA:1242:C:O2'	1:AA:1303:C:H5''	2.17	0.44
22:CV:6167:G:H2'	22:CV:6168:G:O4'	2.18	0.44
1:AA:964:A:H5'	1:AA:1199:U:OP1	2.17	0.44
38:DP:110:ILE:HA	38:DP:110:ILE:HD12	1.85	0.44
39:DQ:60:LEU:HD23	39:DQ:60:LEU:O	2.16	0.44
18:AR:45:SER:HB3	18:AR:51:LEU:CG	2.48	0.44
29:BG:22:GLY:C	29:BG:23:ARG:HD3	2.38	0.44
23:DA:2364:C:H1'	45:DW:36:ILE:HD11	2.00	0.44
19:AS:33:THR:HG22	19:AS:51:VAL:HA	1.99	0.44
38:BP:107:ASP:OD2	38:BP:109:GLU:HB2	2.17	0.44
32:BJ:161:LEU:HD23	32:BJ:161:LEU:N	2.32	0.44
1:CA:833:U:H2'	1:CA:834:C:H6	1.83	0.44
26:BD:93:VAL:HG21	26:BD:177:PRO:HA	1.99	0.44
34:BL:65:ARG:HD2	34:BL:65:ARG:N	2.33	0.44
23:DA:613:U:H4'	23:DA:616:A:N6	2.33	0.44
23:DA:225:A:O2'	23:DA:257:A:H4'	2.17	0.44
1:AA:1468:A:H2'	1:AA:1469:G:O4'	2.18	0.44
23:BA:466:A:N3	23:BA:683:C:H1'	2.32	0.44
29:DG:38:SER:HB2	29:DG:41:MET:HG3	1.99	0.44
23:DA:1668:A:H4'	23:DA:1669:A:O5'	2.18	0.44
27:DE:50:SER:HA	27:DE:92:PRO:O	2.17	0.44
23:DA:2767:C:H2'	23:DA:2768:C:H6	1.83	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
43:DU:63:LYS:HE3	43:DU:63:LYS:HB2	1.81	0.44
23:BA:1733:G:O5'	23:BA:1733:G:H8	2.01	0.44
26:BD:176:ILE:HG22	26:BD:176:ILE:O	2.17	0.44
26:BD:176:ILE:HD12	26:BD:176:ILE:N	2.33	0.44
1:CA:1324:A:O4'	1:CA:136(A):C:H4'	2.18	0.44
51:B3:41:PRO:HG3	51:B3:49:HIS:HE1	1.82	0.44
25:BC:226:MET:HB3	25:BC:230:ASP:HB2	2.00	0.44
2:AB:163:PHE:HA	2:AB:185:ILE:O	2.18	0.44
32:DJ:156:GLN:O	32:DJ:157:ARG:HB2	2.16	0.44
23:DA:2285:C:H41	51:D3:25:LYS:NZ	2.16	0.44
23:DA:956:G:OP1	35:DM:86:GLY:N	2.48	0.44
3:AC:131:ARG:HH21	5:AE:50:GLU:HG2	1.83	0.44
43:BU:75:ILE:HG12	43:BU:76:CYS:N	2.32	0.44
43:BU:81:LYS:NZ	43:BU:99:CYS:SG	2.91	0.44
35:BM:141:GLN:N	44:BV:53:ILE:HB	2.33	0.44
25:BC:35:LYS:HB3	25:BC:36:PRO:HD3	2.00	0.44
47:DY:16:LEU:N	47:DY:16:LEU:HD22	2.33	0.44
48:DZ:46:ASN:HA	48:DZ:46:ASN:HD22	1.66	0.44
23:DA:2543:G:H2'	23:DA:2544:G:O4'	2.17	0.44
25:DC:142:VAL:HG23	25:DC:192:THR:C	2.38	0.44
23:DA:794:G:C5	23:DA:795:C:C4	3.06	0.44
4:AD:3:ARG:NH2	4:AD:118:ARG:HD3	2.30	0.44
47:DY:46:GLN:O	47:DY:47:ASN:CB	2.60	0.44
23:DA:186:G:H2'	23:DA:187:G:H8	1.82	0.44
8:AH:89:PRO:HA	8:AH:92:ARG:NH1	2.33	0.44
20:AT:48:LYS:HD3	20:AT:51:GLU:OE2	2.17	0.44
20:AT:50:GLU:HG3	20:AT:51:GLU:N	2.33	0.44
25:BC:235:GLY:O	25:BC:237:GLU:N	2.45	0.44
1:AA:1298:C:N4	7:AG:114:ARG:HD3	2.33	0.44
26:BD:47:VAL:HG23	26:BD:84:PHE:O	2.18	0.44
25:BC:10:THR:HG23	25:BC:13:ARG:HB2	1.97	0.44
1:CA:754:C:C3'	1:CA:754:C:O2	2.66	0.44
23:DA:2850:A:H8	23:DA:2850:A:H5'	1.80	0.44
2:AB:178:ARG:HB2	2:AB:178:ARG:HH11	1.82	0.44
23:DA:2666:C:H3'	23:DA:2667:C:H6	1.82	0.44
1:AA:1528:U:O2'	1:AA:1530:G:H5'	2.18	0.44
23:BA:813:U:H2'	23:BA:814:C:C6	2.52	0.44
23:BA:2346:A:C2	23:BA:2383:G:C2	3.06	0.44
1:AA:1499:A:H1'	1:AA:1520:G:H5'	2.00	0.44
28:DF:126:ASP:OD2	28:DF:130:ASN:HB2	2.18	0.44
32:BJ:149:PRO:O	32:BJ:150:ASP:HB2	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1071:C:H5''	5:AE:49:PRO:HG2	2.00	0.44
44:BV:5:LEU:HD21	44:BV:39:VAL:HB	1.99	0.44
44:BV:110:GLY:HA2	44:BV:146:ILE:HG23	1.99	0.44
23:BA:1028:A:H61	23:BA:1125:G:H2'	1.83	0.44
1:AA:164:U:H2'	1:AA:165:C:C5	2.53	0.44
36:BN:109:ALA:HA	36:BN:110:PRO:HD3	1.84	0.44
30:DH:25:TYR:O	30:DH:29:TYR:HB3	2.18	0.44
32:DJ:161:LEU:HD23	32:DJ:161:LEU:N	2.32	0.44
13:AM:81:LEU:HD11	13:AM:88:ARG:HH21	1.82	0.44
8:AH:1:MET:HG2	8:AH:2:LEU:O	2.17	0.44
44:BV:54:HIS:HB3	44:BV:101:PRO:HD3	1.99	0.44
1:CA:934:C:H5	1:CA:1344:C:H2'	1.82	0.44
1:AA:1010:G:H2'	1:AA:1011:G:C8	2.53	0.44
26:DD:21:VAL:HG12	26:DD:23:VAL:HG13	1.98	0.44
1:CA:1272:G:H2'	1:CA:1273:G:C8	2.52	0.44
23:DA:2580:U:H4'	26:DD:130:GLY:HA2	1.99	0.44
26:BD:190:GLY:HA2	26:BD:191:PRO:HD3	1.74	0.44
23:BA:1483:G:H2'	23:BA:1484:G:C8	2.53	0.44
27:BE:82:ILE:O	27:BE:82:ILE:HG13	2.17	0.44
23:BA:2450:A:O2'	23:BA:2451:A:H5'	2.18	0.44
23:DA:212:G:O2'	23:DA:213:A:H5'	2.18	0.44
40:DR:27:ALA:HB3	40:DR:61:VAL:HG11	1.99	0.44
39:BQ:92:ARG:CD	39:BQ:95:LEU:H	2.31	0.44
39:BQ:112:ARG:NH2	40:BR:46:VAL:HG21	2.33	0.44
42:BT:21:PHE:CD2	42:BT:26:TYR:HD2	2.35	0.44
1:CA:735:C:C2	1:CA:736:C:C5	3.05	0.44
23:BA:637:A:OP2	34:BL:115:LEU:HB2	2.17	0.44
43:BU:4:LYS:HD3	43:BU:4:LYS:H	1.83	0.44
46:BX:11:ARG:NH1	46:BX:61:ARG:H	2.16	0.44
25:BC:120:GLY:O	25:BC:131:LEU:HB3	2.17	0.44
25:DC:120:GLY:O	25:DC:131:LEU:HB3	2.17	0.44
1:AA:691:G:H3'	11:AK:26:ASN:HD21	1.82	0.44
1:CA:942:G:H2'	1:CA:943:U:C6	2.52	0.44
1:AA:922:G:C6	1:AA:923:A:C6	3.05	0.44
46:DX:43:TYR:HA	46:DX:44:PRO:HD3	1.77	0.44
34:BL:12:ALA:HB1	34:BL:16:ARG:HB2	1.99	0.44
25:DC:69:ARG:HH12	25:DC:117:VAL:HG21	1.82	0.44
46:DX:27:GLU:HG2	46:DX:28:GLY:N	2.31	0.44
23:DA:2376:A:H61	37:DO:89:ARG:HG3	1.83	0.44
23:DA:464:U:H4'	52:D4:5:TRP:CZ3	2.53	0.44
4:AD:172:PRO:HD2	4:AD:173:TRP:CE3	2.53	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:AD:188:LEU:HA	4:AD:189:PRO:HD3	1.77	0.44
40:DR:72:VAL:HG22	40:DR:85:LYS:O	2.18	0.44
12:CL:84:ILE:HG23	12:CL:97:TYR:HB3	2.00	0.44
46:BX:90:ILE:O	46:BX:94:LEU:N	2.50	0.44
1:AA:1187:G:H2'	1:AA:1188:A:H8	1.83	0.44
26:DD:47:VAL:HG23	26:DD:84:PHE:O	2.17	0.44
25:BC:16:MET:HE1	25:BC:208:LYS:HG2	1.99	0.44
23:BA:1308:A:N6	23:BA:1309:G:C2	2.86	0.44
36:BN:11:ASN:O	36:BN:12:ARG:HB2	2.18	0.44
1:AA:505:G:C6	1:AA:535:A:C2	3.06	0.44
1:CA:1499:A:H1'	1:CA:1520:G:H5'	1.99	0.44
1:AA:1528:U:H6	1:AA:1528:U:C5'	2.31	0.44
47:DY:24:LEU:O	47:DY:24:LEU:HD23	2.18	0.44
2:AB:115:LEU:HD12	2:AB:118:LEU:HD12	1.99	0.44
23:DA:534:U:O2'	39:DQ:49:HIS:CD2	2.69	0.44
23:DA:2647:U:H2'	23:DA:2648:C:H6	1.82	0.44
1:CA:1235:U:H5''	21:CU:3:LYS:CD	2.47	0.44
1:AA:554:C:H2'	1:AA:555:C:C6	2.53	0.44
44:DV:5:LEU:HD21	44:DV:39:VAL:HB	2.00	0.44
23:DA:1833:U:C2	23:DA:1834:U:C5	3.06	0.44
23:DA:1655:A:O2'	26:DD:115:GLY:HA2	2.18	0.44
3:CC:33:LEU:HD21	14:CN:53:LEU:CD2	2.46	0.44
44:DV:92:SER:HB2	44:DV:94:GLU:OE2	2.17	0.44
2:CB:131:PRO:O	2:CB:135:GLN:HG3	2.17	0.44
38:DP:107:ASP:OD2	38:DP:109:GLU:HB2	2.18	0.44
28:BF:16:ARG:HB3	28:BF:17:PRO:CD	2.48	0.44
26:DD:77:ILE:HG22	26:DD:78:LEU:N	2.33	0.44
23:BA:649:G:C5	23:BA:650:C:C4	3.05	0.44
9:CI:45:ALA:O	9:CI:48:GLU:HB2	2.17	0.44
7:AG:86:GLN:HB2	7:AG:148:ASN:ND2	2.33	0.44
1:CA:336:C:H2'	1:CA:337:C:C6	2.53	0.44
1:CA:551:U:O2'	12:CL:85:ARG:HD2	2.18	0.44
48:DZ:16:PRO:HB2	48:DZ:18:ASP:OD1	2.18	0.44
23:DA:238:C:O2'	23:DA:608:A:H1'	2.17	0.44
1:CA:358:U:H2'	1:CA:359:U:C6	2.52	0.44
3:AC:44:GLU:OE1	3:AC:52:LEU:HD21	2.18	0.44
1:CA:645:C:H2'	1:CA:646:U:O4'	2.18	0.44
1:AA:1235:U:H5''	21:AU:3:LYS:HB2	2.00	0.44
23:DA:994:C:OP1	39:DQ:53:ARG:NH2	2.51	0.44
32:BJ:78:VAL:O	32:BJ:79:ASN:HB2	2.18	0.44
1:CA:1417:G:N2	1:CA:1482:G:H2'	2.33	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:BL:35:HIS:O	34:BL:36:LYS:HB2	2.17	0.44
1:CA:1399:C:H4'	1:CA:1400:C:C5'	2.19	0.44
39:DQ:79:PHE:HE2	39:DQ:106:PHE:CZ	2.36	0.44
40:DR:7:THR:HG23	40:DR:22:VAL:HG11	1.99	0.44
25:BC:222:ARG:HH12	25:BC:239:ARG:CZ	2.31	0.44
2:CB:163:PHE:HA	2:CB:185:ILE:O	2.18	0.44
28:DF:5:LEU:O	28:DF:8:LYS:HB3	2.18	0.44
42:DT:35:THR:HB	42:DT:38:GLU:H	1.82	0.44
44:DV:163:LEU:H	44:DV:163:LEU:CD2	2.29	0.44
38:BP:27:THR:HG22	38:BP:90:GLN:HB3	1.98	0.44
1:AA:675:A:H2'	1:AA:676:A:C8	2.52	0.44
25:DC:35:LYS:O	25:DC:64:ILE:HD12	2.18	0.44
25:DC:33:LEU:O	25:DC:36:PRO:HD2	2.18	0.44
25:DC:83:GLU:OE1	25:DC:104:TYR:OH	2.31	0.44
2:AB:98:LEU:O	2:AB:101:MET:HG3	2.17	0.44
23:DA:528:A:H2	23:DA:2043:C:C4'	2.28	0.44
23:BA:1652:A:H2'	23:BA:1653:G:O4'	2.18	0.44
23:DA:557:U:H2'	23:DA:558:G:C8	2.52	0.44
1:CA:1511:G:C6	1:CA:1512:U:C4	3.06	0.44
14:AN:52:GLN:O	14:AN:54:PRO:HD3	2.18	0.44
14:CN:52:GLN:O	14:CN:54:PRO:HD3	2.18	0.44
20:AT:72:LEU:C	20:AT:72:LEU:HD23	2.38	0.44
23:BA:598:G:H5'	34:BL:15:ARG:CB	2.48	0.44
6:AF:63:TYR:O	6:AF:65:VAL:HG13	2.17	0.44
32:DJ:62:ARG:CZ	32:DJ:64:ASP:HB2	2.47	0.44
23:DA:301:G:C6	23:DA:317:G:C6	3.06	0.44
13:CM:115:LYS:HB2	13:CM:115:LYS:HE3	1.87	0.44
47:DY:19:VAL:HG12	47:DY:23:LYS:HE3	2.00	0.44
22:AV:6167:G:H2'	22:AV:6168:G:O4'	2.17	0.44
1:CA:622:A:C8	1:CA:623:C:C5	3.05	0.44
41:BS:10:VAL:O	41:BS:12:ILE:N	2.50	0.44
44:DV:110:GLY:HA2	44:DV:146:ILE:HG23	1.99	0.44
44:DV:118:GLN:HG3	44:DV:175:VAL:HG13	1.99	0.44
5:CE:11:ILE:HG13	5:CE:31:LEU:HD22	2.00	0.44
52:D4:24:THR:HG23	52:D4:27:GLY:HA3	1.99	0.44
23:BA:1615:C:C5	23:BA:1617:C:C4	3.06	0.44
26:DD:176:ILE:N	26:DD:176:ILE:HD12	2.33	0.44
23:BA:2591:C:H2'	23:BA:2592:G:C8	2.53	0.44
23:DA:2722:G:H2'	23:DA:2723:C:C6	2.53	0.44
7:CG:86:GLN:HB2	7:CG:148:ASN:ND2	2.33	0.44
23:DA:2768:C:C4	23:DA:2769:C:C5	3.06	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DA:851:U:O2'	48:DZ:42:ALA:O	2.34	0.44
11:CK:38:ASN:HA	11:CK:39:PRO:HD3	1.88	0.44
7:CG:107:ALA:HB2	7:CG:134:ALA:HB2	2.00	0.44
8:AH:13:ILE:O	8:AH:17:THR:HG23	2.17	0.44
6:CF:98:LEU:HD12	6:CF:98:LEU:O	2.18	0.44
23:BA:1526:G:H2'	23:BA:1527:G:C8	2.53	0.44
40:BR:38:LEU:C	40:BR:39:LEU:HD22	2.38	0.44
9:CI:19:LEU:CD2	9:CI:59:PHE:HB3	2.29	0.44
23:BA:2285:C:H41	51:B3:25:LYS:NZ	2.15	0.44
1:CA:1378:C:H5	1:CA:1379:G:C8	2.36	0.44
23:DA:637:A:C5'	34:DL:116:GLY:HA2	2.48	0.44
34:DL:85:LEU:HD21	34:DL:116:GLY:O	2.17	0.44
25:DC:132:PRO:HA	25:DC:190:TYR:HA	2.00	0.44
25:DC:271:ILE:O	25:DC:272:ALA:CB	2.65	0.44
25:BC:33:LEU:O	25:BC:36:PRO:HD2	2.18	0.44
47:BY:16:LEU:N	47:BY:16:LEU:HD22	2.33	0.44
25:BC:228:PRO:HD3	25:BC:234:GLY:O	2.18	0.44
38:DP:74:ARG:HD3	38:DP:76:PHE:CE2	2.53	0.44
36:BN:9:LYS:HD3	36:BN:43:GLU:OE1	2.18	0.44
10:CJ:74:ILE:HG12	10:CJ:74:ILE:O	2.18	0.44
23:BA:2376:A:H61	37:BO:89:ARG:HG3	1.82	0.44
37:DO:26:LEU:HG	37:DO:39:ILE:HD11	2.00	0.44
13:AM:14:ARG:HG2	13:AM:44:ARG:NH1	2.31	0.44
44:BV:141:VAL:HA	44:BV:144:LEU:HD23	2.00	0.44
24:DB:79:C:O5'	24:DB:79:C:H6	2.01	0.44
3:CC:36:ASP:HA	3:CC:39:ILE:HD12	2.00	0.44
4:CD:118:ARG:O	4:CD:122:ARG:HB2	2.18	0.44
24:DB:65:C:O2'	24:DB:66:A:H5'	2.18	0.44
9:AI:10:ARG:HD3	9:AI:11:LYS:N	2.33	0.44
32:DJ:116:THR:HG23	32:DJ:117:HIS:N	2.33	0.44
41:BS:14:PRO:O	41:BS:18:ARG:HG3	2.18	0.44
24:BB:102:G:H1'	44:BV:73:GLN:HE22	1.82	0.44
1:AA:920:U:O4'	1:AA:1080:A:C2	2.71	0.44
23:BA:322:A:P	27:BE:169:ASN:HB2	2.58	0.44
9:AI:77:ILE:O	9:AI:81:ILE:HG13	2.18	0.44
9:CI:18:PHE:HB2	9:CI:62:TYR:HD2	1.83	0.44
6:CF:61:LEU:HB3	6:CF:63:TYR:CE2	2.51	0.44
12:AL:24:PRO:HD2	12:AL:97:TYR:OH	2.17	0.44
1:AA:801:U:H2'	1:AA:802:A:H8	1.81	0.44
23:DA:2562:U:H1'	33:DK:23:ARG:HH12	1.81	0.44
23:DA:2850:A:H5''	23:DA:2868:A:C2	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
12:CL:54:VAL:HG12	12:CL:55:ALA:N	2.32	0.44
23:DA:2809:A:C6	23:DA:2892:A:C8	3.05	0.44
27:BE:11:VAL:HG13	27:BE:196:LEU:HD21	2.00	0.44
23:DA:826:U:H2'	23:DA:828:U:O4'	2.18	0.44
1:AA:270:A:C6	1:AA:271:C:C4	3.05	0.44
25:BC:72:LYS:HE3	25:BC:101:GLU:HB3	1.99	0.44
23:BA:1536:A:H5''	23:BA:1537:C:OP2	2.17	0.44
37:DO:36:TYR:H	37:DO:36:TYR:HD1	1.64	0.44
23:DA:855:G:H2'	23:DA:856:C:H6	1.83	0.44
23:BA:405:U:H3'	23:BA:406:G:C5'	2.48	0.44
26:BD:77:ILE:HG22	26:BD:78:LEU:N	2.33	0.44
13:CM:81:LEU:HD11	13:CM:88:ARG:HH21	1.82	0.44
23:DA:1798:U:H5''	25:DC:259:THR:O	2.17	0.44
29:DG:38:SER:HB2	29:DG:41:MET:CG	2.48	0.44
23:DA:1717:G:C6	23:DA:1743:G:C6	3.06	0.44
3:CC:44:GLU:OE1	3:CC:52:LEU:HD21	2.18	0.44
52:B4:13:ALA:O	52:B4:17:GLY:HA3	2.17	0.44
3:AC:121:ALA:HB2	3:AC:198:VAL:HG21	1.99	0.44
52:B4:3:ARG:HD3	52:B4:3:ARG:HA	1.75	0.44
17:CQ:38:ARG:N	17:CQ:38:ARG:HD2	2.33	0.44
23:DA:2574:G:H8	23:DA:2574:G:O5'	2.01	0.44
1:CA:965:A:C2	1:CA:969:A:C2	3.06	0.44
23:BA:1328:G:H2'	23:BA:1330:C:C5	2.52	0.44
1:AA:814:A:N7	1:AA:816:A:C4	2.85	0.44
23:DA:1191:G:OP1	34:DL:35:HIS:CD2	2.71	0.43
49:B1:46:ASN:HB2	49:B1:64:LYS:CB	2.47	0.43
7:AG:65:ALA:O	7:AG:69:VAL:HG23	2.18	0.43
36:BN:10:LEU:CB	36:BN:17:ARG:NE	2.74	0.43
34:BL:16:ARG:NH2	34:BL:18:ARG:H	2.15	0.43
30:DH:9:LEU:HB2	30:DH:12:LEU:HB2	2.00	0.43
29:DG:42:ARG:O	29:DG:52:VAL:HA	2.18	0.43
1:AA:1295:G:O2'	13:AM:14:ARG:NH1	2.51	0.43
44:BV:106:GLY:HA3	44:BV:142:SER:HG	1.83	0.43
23:BA:1694:C:C5'	23:BA:1694:C:C6	2.96	0.43
30:DH:15:VAL:HG12	30:DH:16:GLY:N	2.32	0.43
23:DA:2777:G:H4'	23:DA:2778:A:H5'	2.00	0.43
2:CB:178:ARG:HB2	2:CB:178:ARG:HH11	1.83	0.43
23:BA:863:A:OP1	35:BM:21:THR:HB	2.18	0.43
1:CA:805:C:H2'	1:CA:806:C:C6	2.51	0.43
34:BL:27:HIS:CE1	40:BR:83:ARG:HH12	2.35	0.43
23:DA:514:A:H1'	23:DA:581:C:O2'	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1117:G:N2	1:CA:1180:A:H1'	2.33	0.43
6:CF:46:ARG:HH12	18:CR:37:VAL:HG21	1.83	0.43
23:DA:265:A:C8	23:DA:266:G:H1'	2.53	0.43
1:AA:818:G:H1'	1:AA:820:U:H5	1.83	0.43
5:AE:11:ILE:HG13	5:AE:31:LEU:HD22	2.00	0.43
23:BA:2094:G:OP1	30:BH:22:LYS:HD2	2.17	0.43
30:BH:29:TYR:C	30:BH:32:PRO:HD2	2.38	0.43
21:AU:11:GLY:O	21:AU:15:ARG:HG3	2.18	0.43
23:DA:150:C:H2'	23:DA:151:C:H6	1.82	0.43
16:AP:24:ALA:C	16:AP:26:ARG:H	2.21	0.43
1:AA:1411:C:O2'	1:AA:1412:C:H5'	2.18	0.43
23:DA:2839:G:C5	23:DA:2840:C:C4	3.06	0.43
3:AC:8:ILE:HD12	3:AC:16:ARG:HH21	1.83	0.43
23:BA:616:A:C4'	23:BA:617:G:OP1	2.66	0.43
1:AA:1448:C:H2'	1:AA:1449:C:C6	2.53	0.43
1:CA:1206:G:H2'	1:CA:1207:G:O4'	2.18	0.43
1:AA:179:A:H2'	1:AA:180:U:C6	2.53	0.43
27:DE:50:SER:HB2	27:DE:94:PRO:HD3	1.99	0.43
23:BA:1285:G:C5	23:BA:1329:U:C4	3.06	0.43
2:AB:113:HIS:O	2:AB:116:GLU:HG2	2.18	0.43
5:AE:41:VAL:HG11	5:AE:113:ALA:HA	2.00	0.43
34:BL:101:VAL:HG23	34:BL:108:LYS:H	1.83	0.43
23:BA:2709:G:O2'	23:BA:2710:C:H5'	2.18	0.43
43:BU:63:LYS:HB2	43:BU:63:LYS:HE3	1.80	0.43
23:DA:843:G:H5''	23:DA:843:G:C8	2.53	0.43
20:AT:36:LEU:HA	20:AT:36:LEU:HD13	1.86	0.43
1:AA:1340:A:O2'	22:AV:6161:U:H5'	2.17	0.43
52:D4:13:ALA:O	52:D4:17:GLY:HA3	2.18	0.43
23:BA:2096:U:H2'	23:BA:2097:C:C6	2.53	0.43
23:BA:941:A:H4'	34:BL:35:HIS:CD2	2.54	0.43
23:DA:251:A:H5''	34:DL:51:PHE:CE1	2.52	0.43
40:DR:45:THR:O	40:DR:46:VAL:C	2.56	0.43
40:BR:40:LEU:H	40:BR:47:VAL:CG2	2.31	0.43
40:BR:6:LYS:O	40:BR:6:LYS:HG3	2.19	0.43
43:DU:14:LEU:C	43:DU:14:LEU:HD23	2.38	0.43
23:BA:114(B):A:O2'	23:BA:1143:A:H3'	2.17	0.43
53:D5:25:MET:SD	53:D5:47:LYS:HG2	2.58	0.43
35:DM:85:LYS:HG3	35:DM:86:GLY:N	2.32	0.43
43:BU:6:HIS:O	43:BU:7:VAL:O	2.35	0.43
29:DG:149:ARG:HD2	29:DG:164:TYR:HE1	1.82	0.43
29:DG:86:GLU:HG2	29:DG:164:TYR:O	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
38:DP:26:ASP:OD1	38:DP:26:ASP:O	2.36	0.43
23:BA:1211:U:H4'	23:BA:1212:G:OP2	2.18	0.43
3:AC:23:TYR:HB2	10:AJ:93:GLY:O	2.19	0.43
36:DN:8:ARG:CZ	36:DN:43:GLU:HG3	2.48	0.43
32:DJ:57:LEU:CD1	32:DJ:142:ARG:HB2	2.48	0.43
23:DA:1678:G:O5'	23:DA:1678:G:C8	2.62	0.43
23:DA:1317:A:C6	23:DA:1318:C:C4	3.05	0.43
23:DA:1335:U:H2'	23:DA:1336:A:C8	2.53	0.43
48:DZ:40:THR:OG1	48:DZ:43:ILE:HG12	2.19	0.43
44:DV:141:VAL:HA	44:DV:144:LEU:HD23	2.00	0.43
37:BO:26:LEU:HG	37:BO:39:ILE:HD11	2.00	0.43
35:DM:101:ARG:HG3	35:DM:102:VAL:N	2.33	0.43
51:D3:11:LEU:HD13	51:D3:12:GLU:H	1.83	0.43
42:BT:70:LEU:HD23	42:BT:70:LEU:C	2.39	0.43
23:BA:479:A:HO2'	23:BA:481:G:H8	1.64	0.43
2:AB:71:VAL:HG12	2:AB:93:VAL:HB	2.00	0.43
9:AI:24:GLY:O	9:AI:26:VAL:HG23	2.19	0.43
29:BG:107:VAL:HG23	29:BG:109:PHE:CE1	2.52	0.43
43:DU:76:CYS:HB3	43:DU:77:PRO:CD	2.48	0.43
1:CA:920:U:H2'	1:CA:921:U:H6	1.80	0.43
23:BA:500:G:N2	23:BA:502:A:H2'	2.33	0.43
47:BY:19:VAL:HG12	47:BY:23:LYS:HE3	2.00	0.43
1:CA:235:C:H2'	1:CA:236:G:H8	1.80	0.43
1:CA:555:C:H2'	1:CA:556:C:H6	1.82	0.43
1:CA:164:U:H2'	1:CA:165:C:C5	2.53	0.43
16:CP:24:ALA:C	16:CP:26:ARG:H	2.21	0.43
37:BO:36:TYR:HD1	37:BO:36:TYR:H	1.64	0.43
23:BA:2257:U:O2'	23:BA:2258:C:H5'	2.18	0.43
19:AS:58:VAL:O	19:AS:58:VAL:HG23	2.18	0.43
45:DW:56:ASP:O	45:DW:57:PHE:HB2	2.19	0.43
1:AA:934:C:C5	1:AA:1344:C:H2'	2.53	0.43
9:AI:48:GLU:N	9:AI:49:PRO:CD	2.80	0.43
1:CA:1361:G:C6	1:CA:136(A):C:C4	3.06	0.43
8:CH:24:THR:HG22	8:CH:25:ASP:N	2.34	0.43
23:BA:699:A:H4'	23:BA:1634:A:N7	2.33	0.43
44:BV:152:ALA:N	44:BV:169:GLU:O	2.49	0.43
30:BH:115:ALA:HB3	30:BH:129:THR:O	2.18	0.43
25:BC:175:LEU:HD12	25:BC:185:VAL:HG21	2.00	0.43
1:CA:1124:G:H5'	10:CJ:35:SER:HB2	1.99	0.43
23:DA:1483:G:H2'	23:DA:1484:G:C8	2.53	0.43
40:BR:89:GLN:HA	40:BR:90:PRO:HD3	1.81	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:BN:33:ARG:HG3	36:BN:115:GLU:HG3	1.98	0.43
3:CC:121:ALA:HB2	3:CC:198:VAL:HG21	1.99	0.43
26:BD:203:LYS:HD2	26:BD:203:LYS:O	2.18	0.43
26:DD:192:ASN:N	26:DD:192:ASN:HD22	2.16	0.43
46:DX:75:GLU:HA	46:DX:75:GLU:OE1	2.18	0.43
1:AA:1403:C:H6	1:AA:1403:C:O5'	2.02	0.43
1:AA:1272:G:H2'	1:AA:1273:G:C8	2.53	0.43
42:DT:62:LYS:O	42:DT:73:ARG:HB2	2.17	0.43
23:DA:833:U:H5''	34:DL:48:PRO:HB2	1.98	0.43
7:AG:111:ARG:HA	7:AG:112:PRO:HD3	1.82	0.43
39:DQ:90:VAL:O	39:DQ:92:ARG:N	2.52	0.43
23:BA:1826:G:H2'	23:BA:1827:C:H6	1.82	0.43
32:BJ:42:GLU:O	32:BJ:42:GLU:HG3	2.18	0.43
28:BF:66:GLN:NE2	28:BF:94:LEU:HB3	2.33	0.43
28:DF:66:GLN:NE2	28:DF:94:LEU:HB3	2.33	0.43
46:BX:10:LYS:O	46:BX:11:ARG:CG	2.66	0.43
48:BZ:22:ALA:O	48:BZ:26:LEU:HG	2.18	0.43
3:AC:148:GLY:HA3	3:AC:203:PHE:HB3	2.01	0.43
5:AE:43:LEU:HD12	5:AE:109:ILE:HD11	1.99	0.43
2:CB:98:LEU:O	2:CB:101:MET:HG3	2.19	0.43
48:DZ:22:ALA:O	48:DZ:26:LEU:HG	2.18	0.43
3:CC:150:LYS:O	3:CC:200:ALA:HA	2.18	0.43
1:AA:927:G:N1	1:AA:1391:U:C2	2.86	0.43
25:BC:30:GLU:HG3	25:BC:63:ARG:CZ	2.48	0.43
28:DF:84:LYS:HG3	28:DF:85:GLY:H	1.83	0.43
23:BA:662:G:OP1	34:BL:18:ARG:HD2	2.18	0.43
8:AH:86:ILE:HD13	8:AH:86:ILE:HA	1.86	0.43
1:CA:1513:A:H2'	1:CA:1514:C:C6	2.53	0.43
4:AD:173:TRP:NE1	4:AD:189:PRO:HG3	2.33	0.43
32:DJ:116:THR:OG1	32:DJ:117:HIS:N	2.51	0.43
44:BV:137:ILE:N	44:BV:137:ILE:HD12	2.33	0.43
51:B3:11:LEU:HD22	51:B3:11:LEU:HA	1.89	0.43
23:BA:2271:G:H2'	23:BA:2272:U:H6	1.82	0.43
23:DA:598:G:H2'	23:DA:599:G:O4'	2.18	0.43
6:AF:61:LEU:HB3	6:AF:63:TYR:CE2	2.51	0.43
1:AA:622:A:C8	1:AA:623:C:C5	3.06	0.43
23:BA:36:G:C5	23:BA:37:C:C5	3.06	0.43
25:DC:145:VAL:HG12	25:DC:146:GLU:O	2.17	0.43
28:BF:126:ASP:OD2	28:BF:130:ASN:HB2	2.18	0.43
3:AC:57:ILE:HD11	3:AC:66:VAL:HG13	2.00	0.43
23:DA:102:G:H8	23:DA:102:G:H5''	1.82	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
18:CR:56:THR:CB	18:CR:58:LEU:HD13	2.47	0.43
23:DA:1291:C:H2'	23:DA:1292:U:H6	1.81	0.43
3:CC:57:ILE:HD11	3:CC:66:VAL:HG13	1.99	0.43
40:DR:28:GLU:OE1	40:DR:31:ALA:HB2	2.18	0.43
23:BA:1655:A:H4'	26:BD:115:GLY:N	2.33	0.43
23:DA:221:A:N7	23:DA:266:G:C5	2.86	0.43
38:DP:77:PRO:HB2	38:DP:80:SER:HB2	2.01	0.43
1:AA:1179:A:H2'	1:AA:1180:A:O4'	2.18	0.43
23:DA:1655:A:H1'	26:DD:113:PHE:CE2	2.53	0.43
23:DA:1655:A:H4'	26:DD:115:GLY:N	2.34	0.43
23:BA:2050:C:H2'	23:BA:2051:A:C8	2.53	0.43
39:BQ:25:TRP:O	39:BQ:28:ARG:HB3	2.18	0.43
3:CC:33:LEU:HD21	14:CN:53:LEU:HD23	2.00	0.43
25:BC:166:GLN:HB2	25:BC:174:ILE:HG22	2.00	0.43
19:AS:53:ASN:ND2	19:AS:56:GLN:H	2.17	0.43
23:DA:363(C):G:H2'	23:DA:363(D):G:H8	1.84	0.43
8:CH:1:MET:HG2	8:CH:2:LEU:O	2.17	0.43
1:AA:909:A:OP1	12:AL:20:LYS:HD2	2.18	0.43
21:CU:11:GLY:O	21:CU:15:ARG:HG3	2.18	0.43
23:BA:1272:A:O2'	23:BA:1273:U:H5'	2.18	0.43
23:DA:1750:G:H2'	23:DA:1751:C:C6	2.53	0.43
45:DW:14:ARG:HB2	45:DW:14:ARG:HE	1.50	0.43
23:BA:2658:C:H4'	29:BG:158:HIS:CE1	2.53	0.43
1:AA:1464:G:O2'	1:AA:1465:C:H5'	2.18	0.43
23:DA:286:C:C2	23:DA:287:C:C5	3.07	0.43
23:BA:1525:G:H2'	23:BA:1526:G:C8	2.54	0.43
8:AH:100:ILE:HA	8:AH:101:PRO:HD3	1.80	0.43
1:CA:620:C:C2	4:CD:135:LEU:HG	2.54	0.43
23:BA:836:G:H2'	23:BA:837:C:C6	2.53	0.43
9:CI:39:GLY:O	9:CI:40:LEU:HD23	2.18	0.43
1:AA:965:A:C2	1:AA:969:A:C2	3.06	0.43
1:AA:781:A:H4'	1:AA:1522:U:O2'	2.18	0.43
23:BA:685:A:H1'	23:BA:688:U:O4	2.18	0.43
1:AA:424:G:H8	1:AA:424:G:O5'	2.00	0.43
6:AF:98:LEU:HD12	6:AF:98:LEU:O	2.18	0.43
23:DA:708:C:O5'	23:DA:708:C:H6	2.01	0.43
1:CA:945:G:H2'	1:CA:945:G:N3	2.33	0.43
12:CL:63:TYR:HB3	12:CL:64:GLU:H	1.66	0.43
32:DJ:78:VAL:O	32:DJ:79:ASN:HB2	2.18	0.43
23:BA:2550:G:C6	23:BA:2551:C:C4	3.07	0.43
1:AA:902:G:H2'	1:AA:903:G:H8	1.83	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BA:997:G:O2'	23:BA:998:C:H5'	2.18	0.43
1:AA:1130:A:N6	1:AA:1144:G:H21	2.16	0.43
5:AE:78:HIS:HD2	8:AH:104:ARG:HD2	1.83	0.43
5:CE:76:ILE:CG1	5:CE:77:PRO:HD2	2.32	0.43
34:DL:65:ARG:N	34:DL:65:ARG:HD2	2.33	0.43
3:CC:70:VAL:O	3:CC:106:VAL:HG23	2.18	0.43
37:BO:98:VAL:O	37:BO:101:LEU:HB2	2.18	0.43
47:BY:9:GLN:O	47:BY:12:GLU:HB3	2.19	0.43
17:AQ:54:GLY:HA3	17:AQ:82:MET:HE2	2.01	0.43
45:DW:32:ARG:CA	45:DW:35:ASN:HD21	2.32	0.43
48:DZ:26:LEU:HD21	48:DZ:46:ASN:HB3	2.00	0.43
25:DC:102:LYS:C	25:DC:103:ARG:HG2	2.37	0.43
37:DO:18:ILE:HA	37:DO:21:THR:OG1	2.18	0.43
38:BP:61:PHE:CZ	38:BP:76:PHE:HB2	2.54	0.43
27:DE:181:LEU:CD2	27:DE:186:ILE:HD11	2.47	0.43
23:DA:390:A:C5	34:DL:71:VAL:HG21	2.52	0.43
11:AK:87:THR:HA	11:AK:91:ARG:NH2	2.28	0.43
23:DA:777:A:O2'	23:DA:778:G:H5'	2.17	0.43
1:CA:1368:G:O2'	1:CA:1369:C:H5'	2.19	0.43
24:DB:73:A:H3'	24:DB:74:U:C6	2.54	0.43
20:CT:71:THR:HG22	20:CT:72:LEU:N	2.34	0.43
34:DL:8:PRO:C	34:DL:10:PRO:HD3	2.39	0.43
23:BA:479:A:H4'	23:BA:480:A:O5'	2.18	0.43
9:CI:17:VAL:HG22	9:CI:63:ILE:HG13	1.99	0.43
26:DD:36:ARG:HD3	26:DD:85:ASN:HD21	1.84	0.43
2:AB:8:LYS:HA	2:AB:217:ARG:NH1	2.33	0.43
33:DK:47:ILE:HA	33:DK:47:ILE:HD12	1.88	0.43
23:DA:510:C:OP1	23:DA:511:U:OP2	2.36	0.43
11:CK:120:ARG:HA	11:CK:121:PRO:HD3	1.79	0.43
32:DJ:149:PRO:O	32:DJ:150:ASP:HB2	2.19	0.43
35:DM:61:GLY:O	44:DV:177:PRO:HA	2.19	0.43
47:BY:21:LEU:CD2	47:BY:22:GLU:HG3	2.49	0.43
22:AV:6166:U:H2'	22:AV:6167:G:O4'	2.18	0.43
3:CC:186:PHE:CG	3:CC:187:ALA:N	2.86	0.43
23:DA:405:U:H3'	23:DA:406:G:C5'	2.48	0.43
23:BA:1050:A:H2'	23:BA:1051:G:C8	2.54	0.43
23:BA:2590:A:H2'	23:BA:2591:C:C6	2.54	0.43
1:AA:380:G:N2	1:AA:384:G:C5	2.86	0.43
23:BA:869:G:C2	23:BA:870:A:C8	3.06	0.43
1:CA:179:A:H2'	1:CA:180:U:C6	2.53	0.43
23:BA:816:C:H2'	23:BA:817:C:H6	1.83	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:BD:107:THR:O	26:BD:190:GLY:HA2	2.18	0.43
40:DR:61:VAL:HG23	40:DR:61:VAL:O	2.19	0.43
23:BA:1818:U:H2'	25:BC:157:ARG:HG3	2.00	0.43
23:BA:1461:G:O2'	23:BA:1462:C:H5'	2.18	0.43
23:BA:1717:G:C6	23:BA:1743:G:C6	3.07	0.43
3:AC:134:ILE:HD11	3:AC:153:VAL:HG22	2.00	0.43
8:CH:114:THR:HG21	8:CH:119:LEU:HD12	2.00	0.43
7:CG:133:GLY:O	7:CG:137:LYS:HG3	2.18	0.43
23:BA:2515:C:H1'	23:BA:2570:G:N2	2.33	0.43
11:AK:23:ALA:HB1	11:AK:88:GLY:HA3	2.01	0.43
23:BA:969:U:H2'	23:BA:970:C:C6	2.53	0.43
48:DZ:30:ARG:HG2	48:DZ:30:ARG:H	1.63	0.43
23:BA:251:A:H5''	34:BL:51:PHE:CE1	2.53	0.43
34:DL:48:PRO:O	34:DL:49:ARG:C	2.57	0.43
23:BA:1826:G:H2'	23:BA:1827:C:C6	2.53	0.43
23:DA:1826:G:C6	23:DA:1827:C:C4	3.06	0.43
23:DA:1825:A:H2'	23:DA:1826:G:C8	2.54	0.43
5:CE:78:HIS:HD2	8:CH:104:ARG:HD2	1.82	0.43
42:DT:21:PHE:CD2	42:DT:26:TYR:HD2	2.36	0.43
23:BA:2247:A:H2'	23:BA:2248:C:H6	1.83	0.43
23:DA:71:A:OP1	23:DA:72:U:H2'	2.18	0.43
23:DA:142:G:H1'	42:DT:37:THR:HG21	2.01	0.43
23:BA:956:G:OP1	35:BM:86:GLY:N	2.48	0.43
43:DU:4:LYS:H	43:DU:4:LYS:HD3	1.83	0.43
30:DH:72:LEU:HD12	30:DH:140:LEU:CD1	2.48	0.43
9:CI:10:ARG:HD3	9:CI:11:LYS:N	2.33	0.43
37:BO:17:ARG:HG2	37:BO:18:ILE:HD13	2.00	0.43
23:DA:2037:G:H2'	23:DA:2038:G:C8	2.53	0.43
1:CA:391:G:C6	1:CA:392:G:C5	3.07	0.43
19:AS:49:ILE:N	19:AS:49:ILE:HD12	2.30	0.43
4:AD:31:CYS:O	4:AD:32:ALA:HB3	2.18	0.43
50:B2:33:CYS:HB2	50:B2:34:PRO:HD2	2.01	0.43
3:AC:36:ASP:HA	3:AC:39:ILE:HD12	2.00	0.43
29:BG:42:ARG:O	29:BG:52:VAL:HA	2.18	0.43
4:AD:104:VAL:HG11	4:AD:146:ILE:HD12	2.00	0.43
25:BC:142:VAL:HG23	25:BC:192:THR:C	2.38	0.43
51:B3:11:LEU:HD13	51:B3:12:GLU:H	1.83	0.43
20:CT:48:LYS:HD3	20:CT:51:GLU:OE2	2.18	0.43
5:CE:90:VAL:O	5:CE:120:THR:HA	2.17	0.43
26:DD:86:PRO:HB2	26:DD:87:GLU:H	1.42	0.43
29:DG:96:ALA:CB	29:DG:105:LEU:HB3	2.48	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DA:36:G:C5	23:DA:37:C:C5	3.06	0.43
23:DA:2061:G:H5''	23:DA:2503:A:C2	2.53	0.43
23:BA:778:G:C5	23:BA:779:U:C4	3.07	0.43
1:AA:687:A:H2'	1:AA:701:C:H41	1.82	0.43
29:BG:20:ALA:HB1	29:BG:21:PRO:CD	2.47	0.43
1:CA:1295:G:H21	1:CA:1302:U:H3	1.66	0.43
23:BA:1693:U:H1'	25:BC:14:ARG:NH2	2.34	0.43
12:AL:54:VAL:HG12	12:AL:55:ALA:N	2.33	0.43
44:BV:8:TYR:HB2	44:BV:38:TYR:CE2	2.52	0.43
1:AA:1116:C:C3'	1:AA:1117:G:H5''	2.48	0.43
22:CV:6166:U:H2'	22:CV:6167:G:O4'	2.18	0.43
23:DA:2572:A:C8	26:DD:144:ARG:HB3	2.53	0.43
23:BA:719:C:H2'	23:BA:720:C:C6	2.53	0.43
28:DF:16:ARG:HB3	28:DF:17:PRO:CD	2.49	0.43
45:DW:27:GLU:HB2	45:DW:69:PHE:CD1	2.54	0.43
1:CA:1454:G:H2'	1:CA:1455:G:H8	1.81	0.43
1:AA:858:G:O6	1:AA:869:G:C8	2.72	0.43
23:BA:532:A:N1	23:BA:2020:A:H1'	2.34	0.43
27:BE:50:SER:HA	27:BE:92:PRO:O	2.17	0.43
23:BA:1748:G:H2'	23:BA:1749:A:H8	1.84	0.43
1:CA:380:G:N2	1:CA:384:G:C5	2.87	0.43
1:CA:102:G:H2'	1:CA:103:C:H6	1.83	0.43
23:DA:1851:U:H2'	23:DA:1852:C:H6	1.83	0.43
23:DA:2867:G:C5	38:DP:23:ARG:NH1	2.87	0.43
23:DA:1416:G:HO2'	23:DA:1417:C:H5	1.61	0.43
1:AA:833:U:C2	1:AA:834:C:C5	3.07	0.43
16:CP:55:ARG:O	16:CP:58:TYR:HB3	2.19	0.43
1:AA:1029:G:N2	1:AA:103(A):A:N7	2.67	0.43
35:DM:134:ARG:O	35:DM:135:ASP:HB2	2.17	0.43
23:DA:454:A:H4'	23:DA:455:C:OP2	2.19	0.43
1:AA:315:A:H5''	1:AA:317:G:OP2	2.17	0.43
23:DA:1563:G:H2'	23:DA:1564:C:H6	1.83	0.43
38:BP:35:LYS:HB2	38:BP:35:LYS:HE3	1.89	0.43
33:DK:99:PHE:N	33:DK:99:PHE:CD1	2.85	0.43
23:DA:685:A:H1'	23:DA:688:U:O4	2.18	0.43
1:CA:1058:G:H2'	1:CA:1059:C:C6	2.53	0.43
4:AD:128:VAL:HA	4:AD:145:GLU:O	2.17	0.43
40:BR:20:LEU:HD12	40:BR:21:ARG:H	1.82	0.43
40:BR:27:ALA:HB3	40:BR:61:VAL:HG11	1.99	0.43
34:DL:47:ASP:HB3	34:DL:48:PRO:CA	2.49	0.43
23:BA:2015:A:N3	50:B2:2:ALA:N	2.66	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
47:DY:9:GLN:O	47:DY:12:GLU:HB3	2.18	0.43
38:BP:26:ASP:OD1	38:BP:26:ASP:O	2.36	0.43
25:DC:25:THR:HG23	25:DC:27:THR:HG22	2.00	0.43
23:DA:744:G:OP1	26:DD:132:HIS:CB	2.67	0.43
16:AP:20:VAL:HG22	16:AP:21:VAL:N	2.33	0.43
12:AL:69:ILE:HA	12:AL:99:ILE:HG22	2.00	0.43
12:CL:82:VAL:HG22	12:CL:83:LEU:H	1.83	0.43
2:CB:91:PRO:CA	2:CB:154:LEU:HD11	2.42	0.43
43:BU:11:ASP:H	43:BU:27:VAL:CG2	2.32	0.43
3:CC:23:TYR:CD2	3:CC:24:ALA:N	2.87	0.43
36:BN:10:LEU:HD22	36:BN:17:ARG:HD3	2.00	0.43
23:BA:2286:A:C8	23:BA:2287:A:N6	2.87	0.43
46:BX:19:GLN:HG2	46:BX:41:ARG:CB	2.46	0.43
19:AS:27:GLU:HB3	19:AS:28:LYS:H	1.69	0.43
4:AD:117:ALA:O	4:AD:121:VAL:HG23	2.18	0.43
4:AD:188:LEU:HD12	4:AD:188:LEU:N	2.33	0.43
4:CD:3:ARG:NH2	4:CD:118:ARG:HD3	2.30	0.43
24:DB:102:G:H1'	44:DV:73:GLN:HE22	1.83	0.43
23:BA:1478:G:HO2'	23:BA:1558:A:H2	1.63	0.43
34:BL:8:PRO:C	34:BL:10:PRO:HD3	2.39	0.43
23:DA:39:C:H2'	23:DA:40:C:C6	2.53	0.43
9:CI:24:GLY:O	9:CI:26:VAL:HG23	2.18	0.43
43:DU:75:ILE:HG13	43:DU:80:GLY:H	1.83	0.43
25:DC:146:GLU:OE2	25:DC:150:LYS:N	2.51	0.43
36:DN:11:ASN:O	36:DN:12:ARG:HB2	2.18	0.43
23:DA:426:C:C2	23:DA:427:U:C5	3.07	0.43
1:CA:1505:G:H5''	1:CA:1506:U:OP1	2.19	0.43
1:CA:1098:C:C2	1:CA:1099:G:C8	3.06	0.43
49:B1:48:ILE:N	49:B1:48:ILE:HD12	2.34	0.43
1:AA:1505:G:H5''	1:AA:1506:U:OP1	2.19	0.43
23:DA:17:G:H2'	23:DA:18:C:C6	2.53	0.43
23:DA:379:G:N2	46:DX:20:ARG:NH1	2.66	0.43
1:CA:934:C:C5	1:CA:1344:C:H2'	2.54	0.43
1:CA:1272:G:H2'	1:CA:1273:G:H8	1.84	0.43
40:BR:61:VAL:O	40:BR:61:VAL:HG23	2.18	0.43
1:CA:584:G:H2'	1:CA:585:G:C8	2.54	0.43
7:AG:133:GLY:O	7:AG:137:LYS:HG3	2.18	0.43
13:CM:96:LEU:HB3	13:CM:97:PRO:HD2	2.01	0.43
23:DA:1464:C:H2'	23:DA:1465:G:C8	2.53	0.43
1:CA:1385:G:C2	1:CA:1386:G:C8	3.06	0.43
1:CA:838:G:N2	1:CA:849:C:C2	2.87	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:AF:46:ARG:HH12	18:AR:37:VAL:HG21	1.84	0.43
24:DB:45:A:H2'	24:DB:45:A:N3	2.34	0.43
1:CA:596:C:H5'	1:CA:596:C:H6	1.84	0.43
46:BX:75:GLU:OE1	46:BX:75:GLU:HA	2.19	0.43
23:BA:2086:U:H2'	23:BA:2087:G:C8	2.54	0.43
1:CA:495:A:H4'	1:CA:496:A:OP1	2.19	0.43
8:AH:24:THR:HG22	8:AH:25:ASP:N	2.34	0.43
39:DQ:112:ARG:NH2	40:DR:46:VAL:HG21	2.34	0.43
1:CA:1130:A:H61	1:CA:1144:G:H21	1.67	0.43
30:DH:92:VAL:O	30:DH:92:VAL:HG22	2.19	0.43
32:DJ:38:LEU:HD13	32:DJ:39:ILE:N	2.33	0.43
34:BL:86:LYS:HG3	34:BL:87:ASP:N	2.34	0.43
10:AJ:48:THR:HG22	10:AJ:62:HIS:ND1	2.34	0.43
10:CJ:32:ALA:HB2	10:CJ:76:ASN:HB2	2.01	0.43
23:DA:2025:C:C2	23:DA:2026:C:C5	3.07	0.43
25:BC:25:THR:HG23	25:BC:27:THR:HG22	2.01	0.43
53:B5:21:LYS:HA	53:B5:54:GLU:OE2	2.18	0.43
23:DA:1579:A:H2'	23:DA:1580:A:C8	2.54	0.43
12:AL:74:HIS:CD2	12:AL:76:LEU:HB2	2.54	0.43
26:DD:170:LEU:HB3	26:DD:184:VAL:HG12	1.99	0.43
3:AC:23:TYR:CD2	3:AC:24:ALA:N	2.86	0.43
23:BA:644:A:C2	23:BA:646:A:C4	3.07	0.43
35:DM:24:GLY:HA2	35:DM:101:ARG:CA	2.47	0.43
9:AI:18:PHE:HB2	9:AI:62:TYR:HD2	1.83	0.43
9:CI:17:VAL:HG21	9:CI:80:GLY:HA3	2.01	0.43
23:BA:2850:A:H5''	23:BA:2868:A:H2	1.83	0.43
1:CA:1190:G:OP2	3:CC:5:ILE:HG23	2.18	0.43
23:DA:1567:A:C8	25:DC:84:TYR:CE2	3.07	0.43
38:DP:19:LEU:HA	38:DP:20:PRO:HD3	1.74	0.43
38:DP:56:GLY:O	38:DP:59:THR:HG22	2.18	0.43
28:DF:128:ARG:HG2	28:DF:129:GLY:H	1.84	0.43
1:CA:683:G:C6	1:CA:684:A:C5	3.06	0.43
35:BM:104:PHE:HE1	35:BM:125:LEU:HD11	1.83	0.43
49:D1:48:ILE:N	49:D1:48:ILE:HD12	2.33	0.43
12:AL:51:LEU:N	12:AL:51:LEU:HD12	2.34	0.43
23:BA:379:G:N2	46:BX:20:ARG:NH1	2.66	0.43
25:DC:72:LYS:HE3	25:DC:101:GLU:HB3	2.00	0.43
4:AD:137:SER:O	4:AD:138:TYR:O	2.37	0.43
23:BA:2839:G:C5	23:BA:2840:C:C4	3.07	0.43
1:CA:1411:C:O2'	1:CA:1412:C:H5'	2.18	0.43
37:BO:38:GLN:HB3	37:BO:47:THR:HG23	1.99	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DA:2590:A:H2'	23:DA:2591:C:C6	2.54	0.43
23:DA:616:A:H4'	23:DA:617:G:OP1	2.18	0.43
16:CP:43:LYS:HA	16:CP:48:TRP:HB3	2.00	0.43
23:BA:2748:A:C2	23:BA:2757:A:C4	3.06	0.43
23:DA:2450:A:O2'	23:DA:2451:A:H5'	2.18	0.43
23:BA:1464:C:H2'	23:BA:1465:G:C8	2.54	0.43
8:CH:100:ILE:HA	8:CH:101:PRO:HD3	1.80	0.43
35:DM:130:LYS:HG2	35:DM:131:ILE:N	2.33	0.43
23:DA:1620:G:O2'	52:D4:2:LYS:HG2	2.18	0.43
36:DN:55:ALA:HA	36:DN:80:PHE:CE1	2.53	0.43
23:DA:2389:G:H5''	23:DA:2390:U:H5'	2.01	0.43
1:AA:697:U:H2'	1:AA:698:G:H5'	2.01	0.43
44:DV:150:LEU:HD23	44:DV:151:HIS:N	2.33	0.43
23:DA:2678:C:H2'	23:DA:2679:A:H8	1.83	0.43
1:CA:484:G:C8	1:CA:486:U:C2	3.06	0.43
44:BV:40:ASP:OD1	44:BV:42:VAL:HG12	2.19	0.43
5:CE:25:ARG:N	5:CE:25:ARG:HD2	2.32	0.43
44:DV:70:LEU:N	44:DV:70:LEU:HD23	2.34	0.43
48:BZ:16:PRO:HB2	48:BZ:18:ASP:OD1	2.18	0.43
23:BA:519:U:H2'	23:BA:520:G:H8	1.82	0.43
27:BE:70:THR:HG23	27:BE:72:ARG:H	1.83	0.43
34:DL:50:ARG:HB2	53:D5:60:LEU:CD1	2.35	0.43
34:DL:50:ARG:HB3	53:D5:60:LEU:HD21	2.00	0.43
15:CO:36:ILE:HD12	15:CO:63:ARG:NH1	2.34	0.43
28:DF:8:LYS:O	28:DF:12:TYR:HD1	2.02	0.43
45:DW:43:THR:HG22	45:DW:43:THR:O	2.19	0.43
29:BG:86:GLU:CD	29:BG:86:GLU:H	2.22	0.43
44:DV:71:VAL:HA	44:DV:87:ASP:O	2.19	0.43
23:DA:2026:C:C2	23:DA:2027:G:C8	3.07	0.43
30:BH:72:LEU:HD12	30:BH:140:LEU:CD1	2.48	0.43
12:AL:82:VAL:HG23	12:AL:106:ALA:CB	2.44	0.43
7:CG:65:ALA:O	7:CG:69:VAL:HG23	2.18	0.43
36:BN:8:ARG:CZ	36:BN:43:GLU:HG3	2.49	0.43
23:BA:644:A:O2'	23:BA:645:C:H5''	2.18	0.43
37:BO:31:SER:HB3	37:BO:34:HIS:H	1.84	0.43
1:CA:586:C:O2'	1:CA:878:G:H4'	2.19	0.43
23:BA:322:A:O4'	23:BA:340:A:H1'	2.19	0.43
25:BC:235:GLY:C	25:BC:237:GLU:H	2.22	0.43
1:AA:1148:U:H4'	9:AI:14:VAL:HG11	2.01	0.43
2:CB:71:VAL:HG12	2:CB:93:VAL:HB	2.01	0.43
5:CE:92:LYS:O	5:CE:118:ILE:HD12	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DA:214:G:O2'	23:DA:215:G:O4'	2.33	0.43
7:AG:150:ALA:HA	11:AK:59:TYR:HD2	1.84	0.43
1:CA:932:C:H2'	1:CA:933:G:H8	1.83	0.43
4:CD:79:PHE:CD1	4:CD:207:TYR:HD1	2.37	0.43
1:AA:1004:A:H8	1:AA:1026:G:N7	2.17	0.43
37:DO:93:LYS:HE3	37:DO:93:LYS:HA	2.01	0.43
39:DQ:25:TRP:O	39:DQ:28:ARG:HB3	2.19	0.43
51:B3:44:ARG:HB3	51:B3:45:LYS:H	1.67	0.43
44:BV:118:GLN:HG3	44:BV:175:VAL:HG13	2.00	0.43
23:DA:2050:C:H2'	23:DA:2051:A:C8	2.54	0.43
19:AS:51:VAL:HG12	19:AS:52:TYR:O	2.19	0.43
32:BJ:160:LYS:HD2	32:BJ:161:LEU:H	1.83	0.43
4:AD:57:ARG:HB3	4:AD:206:PHE:HB2	2.00	0.43
26:BD:78:LEU:C	26:BD:79:ARG:HD2	2.38	0.43
23:DA:1796:U:H2'	23:DA:1797:C:C6	2.53	0.43
1:CA:1464:G:O2'	1:CA:1465:C:H5'	2.18	0.43
1:CA:774:G:H2'	1:CA:775:G:H5'	2.00	0.43
1:CA:1342:C:O2'	1:CA:1343:G:H5'	2.18	0.43
1:AA:1385:G:C2	1:AA:1386:G:C8	3.07	0.43
8:AH:85:ARG:HH12	8:AH:134:ILE:HG23	1.84	0.43
29:BG:78:GLY:HA2	29:BG:83:TYR:CE1	2.54	0.43
23:BA:708:C:O5'	23:BA:708:C:H6	2.01	0.43
1:CA:424:G:O5'	1:CA:424:G:H8	2.01	0.43
23:BA:1239:G:O2'	23:BA:1240:U:H5'	2.18	0.43
1:AA:290:C:H6	1:AA:290:C:O5'	2.02	0.43
23:BA:843:G:C8	23:BA:843:G:H5''	2.54	0.43
15:AO:65:ARG:O	15:AO:68:ARG:HB2	2.18	0.43
23:DA:2509:G:C5	23:DA:2510:C:C5	3.06	0.43
1:CA:1145:C:H4'	1:CA:1146:A:H8	1.84	0.43
23:DA:587:C:C6	23:DA:671:C:H1'	2.54	0.43
39:DQ:79:PHE:HE1	39:DQ:83:LEU:CD2	2.31	0.43
39:BQ:79:PHE:HE2	39:BQ:106:PHE:CZ	2.36	0.43
25:DC:222:ARG:HE	25:DC:222:ARG:HB2	1.70	0.43
34:BL:114:ILE:CD1	34:BL:130:PHE:CD1	2.92	0.43
43:BU:13:VAL:CG1	43:BU:72:VAL:HB	2.48	0.43
49:D1:64:LYS:HA	49:D1:64:LYS:HE3	2.00	0.43
28:DF:55:LYS:HD2	28:DF:58:GLN:NE2	2.21	0.43
44:BV:71:VAL:HA	44:BV:87:ASP:O	2.18	0.43
25:DC:35:LYS:HB3	25:DC:36:PRO:HD3	2.00	0.43
53:D5:21:LYS:HA	53:D5:54:GLU:OE2	2.19	0.43
36:DN:4:LEU:C	36:DN:6:SER:N	2.72	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:CC:148:GLY:HA3	3:CC:203:PHE:HB3	2.00	0.43
25:BC:35:LYS:HZ1	25:BC:104:TYR:H	1.66	0.43
34:BL:57:THR:CG2	34:BL:59:LEU:HB2	2.48	0.43
23:DA:330:A:O2'	23:DA:331:A:C8	2.72	0.43
28:BF:84:LYS:HG3	28:BF:85:GLY:H	1.84	0.43
23:DA:1678:G:H22	23:DA:1989:G:H22	1.64	0.43
1:AA:265:G:H5'	17:AQ:64:PRO:O	2.19	0.43
4:CD:172:PRO:HD2	4:CD:173:TRP:CE3	2.53	0.43
23:DA:795:C:O2'	23:DA:796:C:H5'	2.18	0.43
4:AD:104:VAL:O	4:AD:108:LEU:HB2	2.19	0.43
9:AI:17:VAL:HG21	9:AI:80:GLY:HA3	2.00	0.43
23:BA:2562:U:H1'	33:BK:23:ARG:HH12	1.81	0.43
1:AA:932:C:H2'	1:AA:933:G:H8	1.84	0.43
23:DA:448:U:H1'	27:DE:84:VAL:HG21	2.01	0.43
23:DA:2563:U:O2	23:DA:2565:A:C8	2.71	0.43
34:DL:105:LEU:H	34:DL:105:LEU:HD12	1.83	0.43
18:CR:84:LYS:HZ2	18:CR:84:LYS:HA	1.82	0.43
44:BV:38:TYR:CD1	44:BV:38:TYR:O	2.72	0.43
23:DA:1028:A:N3	23:DA:2486:G:O2'	2.38	0.43
36:BN:104:ARG:HB2	36:BN:104:ARG:HH11	1.84	0.43
23:DA:1536:A:O5'	23:DA:1536:A:H8	2.02	0.43
23:BA:433:C:O2'	23:BA:434:U:H5'	2.19	0.43
26:DD:93:VAL:HG21	26:DD:177:PRO:HA	1.99	0.43
3:CC:8:ILE:HD12	3:CC:16:ARG:HH21	1.83	0.43
1:CA:1057:G:H2'	1:CA:1058:G:O4'	2.19	0.43
1:CA:782:A:O3'	1:CA:1515:C:H4'	2.19	0.43
7:CG:155:ARG:O	7:CG:156:TRP:CD1	2.72	0.43
1:AA:66:G:H4'	1:AA:173:U:C5	2.53	0.43
23:DA:2086:U:H2'	23:DA:2087:G:C8	2.54	0.43
7:AG:87:VAL:HG11	7:AG:154:TYR:O	2.19	0.43
23:BA:2768:C:C4	23:BA:2769:C:C5	3.07	0.43
1:AA:484:G:C8	1:AA:486:U:C2	3.07	0.43
15:AO:7:GLU:HA	15:AO:10:LYS:HB3	2.01	0.43
23:DA:2585:U:H4'	23:DA:2586:C:OP1	2.17	0.43
23:BA:2330:G:O2'	45:BW:41:ARG:HB2	2.17	0.43
23:DA:2683:C:OP1	38:DP:53:ARG:NH2	2.51	0.43
12:AL:123:LYS:HA	12:AL:124:PRO:HD3	1.89	0.43
7:CG:60:LYS:HA	7:CG:60:LYS:HD2	1.91	0.43
23:BA:2891:G:H8	23:BA:2891:G:O5'	2.01	0.43
1:CA:781:A:H4'	1:CA:1522:U:O2'	2.18	0.43
25:BC:140:THR:O	25:BC:165:ILE:HD12	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BA:1191:G:OP1	34:BL:35:HIS:CD2	2.72	0.43
1:CA:1129:C:H1'	1:CA:1130:A:OP2	2.18	0.43
28:BF:5:LEU:O	28:BF:8:LYS:HB3	2.19	0.43
45:BW:43:THR:O	45:BW:43:THR:HG22	2.19	0.43
29:DG:86:GLU:CD	29:DG:86:GLU:H	2.22	0.43
23:DA:848:G:N9	23:DA:933:A:H8	2.17	0.43
44:DV:51:ALA:HB1	44:DV:57:ILE:HD11	2.01	0.43
48:BZ:26:LEU:HB2	48:BZ:28:LEU:CD1	2.48	0.43
23:BA:643:A:C2	23:BA:644:A:C4	3.07	0.43
23:BA:1335:U:H2'	23:BA:1336:A:C8	2.54	0.43
28:DF:161:THR:HG21	28:DF:172:LEU:HD23	2.01	0.43
23:BA:907:U:O2'	35:BM:101:ARG:NH2	2.46	0.43
4:AD:118:ARG:O	4:AD:122:ARG:HB2	2.18	0.43
24:DB:106:G:O2'	24:DB:107:U:H5'	2.19	0.43
40:DR:72:VAL:CG2	40:DR:85:LYS:HB3	2.48	0.43
24:BB:65:C:O2'	24:BB:66:A:H5'	2.18	0.43
23:BA:2777:G:H4'	23:BA:2778:A:H5'	2.01	0.43
36:DN:2:ARG:HB3	36:DN:3:HIS:CE1	2.54	0.43
42:BT:93:GLU:O	42:BT:94:GLY:C	2.56	0.43
23:BA:102:G:C8	23:BA:102:G:H5''	2.54	0.43
2:CB:115:LEU:HD12	2:CB:118:LEU:HD12	1.99	0.43
33:DK:86:ILE:N	33:DK:86:ILE:HD12	2.31	0.43
23:BA:302:C:O2'	23:BA:303:U:H5'	2.19	0.43
23:DA:1996:C:H5	33:DK:32:TYR:OH	2.02	0.43
22:AV:6164:A:H2'	22:AV:6165:G:C8	2.54	0.43
1:CA:1281:U:H5'	1:CA:1282:C:H5	1.83	0.43
16:CP:50:LYS:HD3	16:CP:51:VAL:N	2.34	0.43
51:D3:30:THR:CG2	51:D3:31:PRO:HD2	2.48	0.43
23:DA:2591:C:H2'	23:DA:2592:G:C8	2.54	0.43
23:BA:855:G:H2'	23:BA:856:C:H6	1.84	0.43
41:DS:70:TYR:H	41:DS:70:TYR:HD2	1.66	0.43
1:CA:337:C:H2'	1:CA:338:A:H8	1.83	0.43
27:DE:117:ARG:HD2	27:DE:190:GLU:O	2.18	0.43
1:AA:757:U:H5''	1:AA:822:C:O2	2.19	0.43
44:DV:128:VAL:CG2	44:DV:132:ASN:HB2	2.49	0.43
25:DC:140:THR:O	25:DC:165:ILE:HD12	2.19	0.43
1:AA:1206:G:H2'	1:AA:1207:G:O4'	2.19	0.43
23:DA:533:G:N3	39:DQ:45:TYR:CE1	2.87	0.43
23:DA:1576:U:C2	23:DA:1577:C:C5	3.07	0.43
23:BA:1232:G:H2'	23:BA:1233:C:H6	1.84	0.43
1:CA:967:C:H4'	9:CI:125:TYR:OH	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
17:AQ:38:ARG:N	17:AQ:38:ARG:HD2	2.33	0.43
4:CD:188:LEU:N	4:CD:188:LEU:HD12	2.33	0.43
47:BY:15:LYS:HE2	47:BY:15:LYS:HA	2.01	0.43
23:DA:944:G:H5'	23:DA:945:A:C5'	2.49	0.43
23:BA:1961:C:O2'	23:BA:1962:C:H5'	2.18	0.43
1:AA:809:G:C6	1:AA:810:C:C5	3.07	0.43
34:DL:101:VAL:HG23	34:DL:108:LYS:H	1.84	0.43
32:DJ:85:VAL:HG22	32:DJ:89:LYS:HG3	2.01	0.43
34:BL:80:TYR:CE1	34:BL:111:ARG:HG2	2.53	0.42
23:DA:783:A:C3'	23:DA:783:A:C8	3.00	0.42
53:B5:32:LEU:HD23	53:B5:33:ASN:H	1.83	0.42
27:BE:63:LYS:HZ1	27:BE:67:GLN:NE2	2.17	0.42
50:D2:33:CYS:HB2	50:D2:34:PRO:HD2	2.01	0.42
46:DX:11:ARG:NH1	46:DX:61:ARG:H	2.17	0.42
25:BC:132:PRO:HA	25:BC:190:TYR:HA	2.01	0.42
44:BV:48:PHE:CE2	44:BV:71:VAL:HG21	2.54	0.42
25:DC:33:LEU:HD23	25:DC:33:LEU:N	2.25	0.42
25:BC:27:THR:O	25:BC:27:THR:CG2	2.62	0.42
25:BC:32:SER:O	25:BC:33:LEU:O	2.36	0.42
43:BU:17:SER:HB2	43:BU:71:LYS:HD2	2.01	0.42
23:DA:2286:A:C8	23:DA:2287:A:N6	2.87	0.42
26:BD:197:ILE:HD11	26:BD:199:ARG:HE	1.84	0.42
23:BA:661:C:H2'	23:BA:662:G:H8	1.83	0.42
28:BF:161:THR:HG21	28:BF:172:LEU:HD23	2.01	0.42
19:AS:29:ARG:HB2	19:AS:48:THR:N	2.33	0.42
30:DH:114:LEU:HA	30:DH:130:TYR:CD1	2.54	0.42
23:DA:775:G:C2	23:DA:777:A:N6	2.86	0.42
50:B2:44:THR:HG22	50:B2:45:VAL:N	2.34	0.42
35:BM:24:GLY:HA2	35:BM:100:GLY:C	2.39	0.42
23:BA:1405:U:H2'	23:BA:1406:U:H6	1.79	0.42
8:CH:89:PRO:HA	8:CH:92:ARG:NH1	2.34	0.42
23:BA:477:A:H2'	23:BA:478:A:C8	2.54	0.42
26:BD:36:ARG:HD3	26:BD:85:ASN:HD21	1.84	0.42
2:AB:11:LEU:HD12	2:AB:217:ARG:NH2	2.34	0.42
29:DG:107:VAL:HG23	29:DG:109:PHE:CE1	2.54	0.42
42:DT:93:GLU:O	42:DT:94:GLY:C	2.57	0.42
15:AO:50:HIS:O	15:AO:53:HIS:HB3	2.19	0.42
23:DA:2243:U:H2'	23:DA:2244:U:H6	1.84	0.42
1:AA:1117:G:N2	1:AA:1180:A:H1'	2.33	0.42
26:BD:111:ARG:HD2	26:BD:160:TYR:HE1	1.84	0.42
23:BA:1576:U:C2	23:BA:1577:C:C5	3.07	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
22:CV:6167:G:H2'	22:CV:6168:G:C8	2.53	0.42
41:DS:10:VAL:HG21	41:DS:103:ILE:HD13	2.01	0.42
8:AH:36:LEU:HD23	8:AH:39:LEU:HD23	2.00	0.42
23:BA:116:C:O2'	23:BA:117:G:H5'	2.19	0.42
23:BA:2846:G:H2'	23:BA:2847:U:C6	2.53	0.42
32:DJ:160:LYS:HD2	32:DJ:161:LEU:H	1.83	0.42
23:DA:2723:C:O5'	23:DA:2723:C:H6	2.02	0.42
53:B5:29:LYS:O	53:B5:29:LYS:HG2	2.19	0.42
32:DJ:27:TYR:O	32:DJ:29:PRO:HD3	2.19	0.42
23:BA:993:G:H1'	40:BR:89:GLN:OE1	2.19	0.42
23:BA:26:G:C6	23:BA:27:G:N1	2.88	0.42
5:CE:41:VAL:HG11	5:CE:113:ALA:HA	2.01	0.42
23:BA:2389:G:H5''	23:BA:2390:U:H5'	2.00	0.42
1:AA:386:C:H2'	1:AA:387:U:H5''	2.00	0.42
25:DC:226:MET:HB3	25:DC:230:ASP:HB2	2.00	0.42
23:BA:460:A:C6	23:BA:470:A:C8	3.07	0.42
39:BQ:59:ARG:O	39:BQ:63:VAL:HG23	2.19	0.42
1:CA:375:U:H4'	16:CP:17:TYR:CE2	2.54	0.42
4:CD:156:GLU:O	4:CD:160:GLN:HG3	2.19	0.42
8:AH:38:ILE:HD12	8:AH:118:VAL:HG12	2.00	0.42
13:AM:96:LEU:HB3	13:AM:97:PRO:HD2	2.00	0.42
23:DA:195:A:H4'	23:DA:251:A:O2'	2.20	0.42
39:BQ:90:VAL:O	39:BQ:92:ARG:N	2.52	0.42
23:BA:2394:C:H2'	23:BA:2395:C:H6	1.84	0.42
42:BT:35:THR:HB	42:BT:38:GLU:H	1.84	0.42
23:DA:142:G:H1'	42:DT:37:THR:CG2	2.49	0.42
28:BF:8:LYS:O	28:BF:12:TYR:HD1	2.02	0.42
37:BO:49:VAL:HG13	37:BO:76:LYS:HB2	2.01	0.42
23:BA:2712:U:O2'	23:BA:2713:A:H5'	2.19	0.42
23:BA:848:G:N9	23:BA:933:A:H8	2.16	0.42
25:DC:31:LYS:HG3	25:DC:33:LEU:HG	2.01	0.42
25:DC:33:LEU:HB2	25:DC:34:VAL:H	1.64	0.42
34:DL:57:THR:C	34:DL:59:LEU:H	2.21	0.42
34:DL:6:LEU:CD2	34:DL:6:LEU:H	2.26	0.42
53:B5:14:VAL:HG22	53:B5:24:ALA:HB2	2.01	0.42
23:BA:1310:G:C3'	23:BA:1311:G:H5''	2.49	0.42
37:DO:100:ALA:HA	37:DO:103:GLU:HB3	2.00	0.42
23:DA:141(A):A:C8	23:DA:1408:C:H1'	2.54	0.42
36:DN:10:LEU:HD22	36:DN:17:ARG:HD3	2.01	0.42
23:BA:2343:C:O2'	23:BA:2344:U:H5'	2.19	0.42
27:DE:139:PHE:HB2	27:DE:166:ALA:CB	2.48	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BA:1612:C:O3'	52:B4:5:TRP:HD1	2.02	0.42
11:CK:91:ARG:HD3	18:CR:88:LYS:HE2	2.00	0.42
23:DA:66:C:H2'	23:DA:67:U:H6	1.84	0.42
24:BB:73:A:C4	24:BB:104:A:C2	3.07	0.42
23:DA:811:U:H3'	34:DL:25:SER:O	2.19	0.42
38:BP:98:LYS:HB3	38:BP:100:TYR:HE1	1.82	0.42
5:CE:92:LYS:HB3	5:CE:119:LEU:HB2	2.01	0.42
28:BF:129:GLY:HA3	28:BF:163:ALA:HB3	2.01	0.42
23:DA:500:G:N2	23:DA:502:A:H2'	2.34	0.42
1:AA:1098:C:C2	1:AA:1099:G:C8	3.07	0.42
10:AJ:6:ILE:HB	10:AJ:98:ILE:HG12	2.01	0.42
2:AB:112:VAL:O	2:AB:115:LEU:HB3	2.18	0.42
38:BP:64:ARG:HA	38:BP:72:VAL:O	2.19	0.42
1:CA:1179:A:H2'	1:CA:1180:A:O4'	2.18	0.42
16:AP:50:LYS:HD3	16:AP:51:VAL:N	2.34	0.42
6:AF:75:LEU:C	6:AF:75:LEU:HD23	2.39	0.42
22:CV:6164:A:H2'	22:CV:6165:G:C8	2.54	0.42
46:DX:45:ASN:HD21	46:DX:47:GLN:HE21	1.67	0.42
30:BH:66:GLU:O	30:BH:70:GLU:HG2	2.20	0.42
41:DS:84:ARG:HB2	41:DS:96:ILE:CG2	2.49	0.42
46:BX:45:ASN:HD21	46:BX:47:GLN:NE2	2.17	0.42
23:BA:2852:G:H2'	23:BA:2853:C:C6	2.54	0.42
23:BA:2846:G:H2'	23:BA:2847:U:H6	1.84	0.42
39:BQ:60:LEU:HD23	39:BQ:60:LEU:O	2.19	0.42
23:DA:816:C:H2'	23:DA:817:C:H6	1.84	0.42
23:BA:616:A:H4'	23:BA:617:G:OP1	2.19	0.42
23:BA:2626:C:H2'	23:BA:2627:G:O4'	2.19	0.42
23:DA:1599:C:H2'	23:DA:1600:C:H6	1.84	0.42
1:AA:833:U:H2'	1:AA:834:C:H6	1.84	0.42
1:AA:37:U:OP1	12:AL:122:LYS:HG3	2.19	0.42
25:DC:202:LYS:HG3	25:DC:203:ASN:OD1	2.19	0.42
1:AA:246:A:C2	1:AA:282:A:C5	3.07	0.42
1:AA:152:A:N6	1:AA:170:U:C2	2.87	0.42
1:AA:156:G:C2	1:AA:166:G:C2	3.07	0.42
23:BA:284:U:H2'	23:BA:285:C:C6	2.54	0.42
29:BG:122:THR:O	29:BG:134:SER:HB2	2.19	0.42
23:DA:1526:G:H2'	23:DA:1527:G:C8	2.54	0.42
17:CQ:7:THR:HG22	17:CQ:58:GLU:HG2	2.00	0.42
32:BJ:93:LYS:HB3	32:BJ:110:LEU:HB2	2.01	0.42
23:DA:735:A:H3'	23:DA:736:C:C6	2.53	0.42
23:BA:1688:U:O2	23:BA:1700:A:H8	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:66:G:H4'	1:CA:173:U:C5	2.54	0.42
26:DD:110:GLY:HA2	26:DD:162:ALA:HB2	2.02	0.42
23:BA:2527:C:O5'	23:BA:2527:C:H6	2.01	0.42
1:AA:1342:C:O2'	1:AA:1343:G:H5'	2.18	0.42
23:DA:1581:G:H2'	23:DA:1582:C:O4'	2.19	0.42
23:BA:1546:A:H2'	23:BA:154(B):C:O4'	2.19	0.42
23:DA:993:G:H1'	40:DR:89:GLN:OE1	2.19	0.42
23:DA:1683:C:H2'	23:DA:1684:C:C6	2.54	0.42
1:AA:1130:A:H61	1:AA:1144:G:H21	1.67	0.42
40:DR:40:LEU:H	40:DR:47:VAL:CG2	2.33	0.42
39:BQ:92:ARG:HH21	40:BR:11:GLN:H	1.67	0.42
15:CO:33:THR:HA	15:CO:63:ARG:HH12	1.83	0.42
32:DJ:40:ASP:OD1	32:DJ:42:GLU:HG2	2.20	0.42
27:BE:63:LYS:NZ	27:BE:67:GLN:NE2	2.67	0.42
23:BA:910:A:C6	23:BA:911:A:C6	3.08	0.42
26:DD:91:VAL:CB	26:DD:95:ILE:HD11	2.42	0.42
35:DM:141:GLN:OXT	44:DV:98:MET:HE3	2.19	0.42
35:BM:141:GLN:NE2	44:BV:72:ARG:HG2	2.35	0.42
53:D5:53:PRO:C	53:D5:57:ARG:NH2	2.73	0.42
9:CI:14:VAL:HG12	9:CI:15:ALA:N	2.34	0.42
26:BD:171:GLU:O	26:BD:171:GLU:HG3	2.19	0.42
2:CB:80:ILE:HD12	2:CB:211:ILE:HB	2.00	0.42
23:DA:2037:G:C6	23:DA:2038:G:C6	3.06	0.42
33:BK:71:ARG:HH12	38:BP:74:ARG:NH2	2.11	0.42
23:DA:2287:A:C2	23:DA:2289:G:C8	3.07	0.42
28:BF:13:GLU:O	28:BF:14:GLU:HB2	2.19	0.42
11:AK:91:ARG:HD3	18:AR:88:LYS:HE2	2.01	0.42
4:CD:122:ARG:O	4:CD:134:ASP:HB2	2.19	0.42
44:BV:102:LEU:HD21	44:BV:124:ILE:HD11	2.02	0.42
1:CA:668:G:H4'	15:CO:48:LYS:HB2	2.02	0.42
23:BA:94:G:N3	47:BY:47:ASN:ND2	2.66	0.42
20:CT:50:GLU:HG3	20:CT:51:GLU:N	2.33	0.42
23:BA:2272:U:H5''	23:BA:2273:A:OP1	2.19	0.42
23:DA:598:G:H5'	34:DL:15:ARG:CG	2.50	0.42
33:DK:7:TYR:CE1	33:DK:20:MET:HB2	2.54	0.42
1:AA:625:G:H2'	1:AA:626:U:C6	2.54	0.42
42:DT:70:LEU:HD23	42:DT:70:LEU:C	2.39	0.42
1:CA:1528:U:C6	1:CA:1528:U:H5''	2.49	0.42
1:CA:1528:U:O2'	1:CA:1530:G:H5'	2.19	0.42
23:BA:840:C:H2'	23:BA:841:A:C8	2.54	0.42
23:BA:2563:U:H4'	33:BK:28:SER:HA	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
38:BP:56:GLY:O	38:BP:59:THR:HG22	2.20	0.42
23:BA:628:G:C6	23:BA:636:G:C2	3.07	0.42
1:AA:1191:A:H5''	3:AC:4:LYS:HZ3	1.85	0.42
36:DN:104:ARG:HH11	36:DN:104:ARG:HB2	1.85	0.42
4:CD:137:SER:O	4:CD:138:TYR:O	2.37	0.42
23:BA:1790:C:H5''	23:BA:1791:A:P	2.59	0.42
8:CH:36:LEU:HD23	8:CH:39:LEU:HD23	2.00	0.42
23:BA:1536:A:O5'	23:BA:1536:A:H8	2.02	0.42
23:DA:1344:G:H5'	23:DA:1384:A:C6	2.55	0.42
1:CA:927:G:N1	1:CA:1391:U:C2	2.87	0.42
23:DA:755:C:H2'	23:DA:756:C:C6	2.54	0.42
4:AD:51:PRO:HB3	4:AD:55:ALA:CB	2.49	0.42
23:DA:909:A:C4	23:DA:912:C:C5	3.07	0.42
23:BA:471:A:H8	23:BA:471:A:O5'	2.02	0.42
11:AK:41:THR:HG21	11:AK:71:LYS:HB2	2.01	0.42
7:AG:95:ARG:O	7:AG:99:LEU:HG	2.19	0.42
23:BA:1683:C:H2'	23:BA:1684:C:C6	2.55	0.42
1:CA:338:A:C6	1:CA:339:C:N4	2.87	0.42
1:AA:1207:G:H2'	1:AA:1208:C:H6	1.83	0.42
44:BV:128:VAL:CG2	44:BV:132:ASN:HB2	2.49	0.42
1:CA:152:A:N6	1:CA:170:U:C2	2.87	0.42
23:BA:286:C:H2'	23:BA:287:C:H6	1.83	0.42
8:AH:84:ARG:O	8:AH:135:CYS:HB2	2.20	0.42
15:AO:8:LYS:O	15:AO:12:ILE:HG13	2.20	0.42
23:BA:1414:G:C4	23:BA:1415:U:C5	3.07	0.42
1:CA:757:U:H5''	1:CA:822:C:O2	2.19	0.42
23:DA:571:A:C8	23:DA:2030:A:N6	2.87	0.42
23:DA:1831:G:C5	23:DA:1832:C:C5	3.07	0.42
23:BA:2364:C:H1'	45:BW:36:ILE:HD11	2.01	0.42
6:AF:9:VAL:HG13	6:AF:59:TYR:O	2.19	0.42
32:BJ:85:VAL:HG22	32:BJ:89:LYS:HG3	2.00	0.42
23:DA:2891:G:O5'	23:DA:2891:G:H8	2.02	0.42
47:DY:15:LYS:HE2	47:DY:15:LYS:HA	2.00	0.42
36:BN:55:ALA:HA	36:BN:80:PHE:CE1	2.54	0.42
24:DB:16:G:C6	24:DB:69:G:C2	3.07	0.42
1:AA:1129:C:H1'	1:AA:1130:A:OP2	2.18	0.42
5:AE:76:ILE:HG12	5:AE:77:PRO:CD	2.31	0.42
43:DU:6:HIS:O	43:DU:7:VAL:O	2.37	0.42
32:BJ:157:ARG:O	32:BJ:157:ARG:HG2	2.20	0.42
32:BJ:38:LEU:HD13	32:BJ:39:ILE:N	2.34	0.42
23:DA:114(B):A:O2'	23:DA:1143:A:H3'	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
42:BT:30:VAL:HG21	42:BT:79:ALA:CB	2.50	0.42
1:CA:979:C:H42	14:CN:18:VAL:CG1	2.28	0.42
23:BA:637:A:C5'	34:BL:116:GLY:HA2	2.49	0.42
23:DA:910:A:C6	23:DA:911:A:C6	3.08	0.42
3:CC:131:ARG:HH21	5:CE:50:GLU:CG	2.32	0.42
3:AC:110:ASN:O	3:AC:141:VAL:HG22	2.19	0.42
27:DE:65:TRP:HZ3	27:DE:75:HIS:HD2	1.60	0.42
23:BA:572:A:H2'	23:BA:573:G:O4'	2.20	0.42
48:DZ:50:VAL:O	48:DZ:54:VAL:HG22	2.19	0.42
23:DA:1006:C:C2	23:DA:1138:G:N2	2.87	0.42
23:DA:1493:C:C4	23:DA:2210:G:O2'	2.73	0.42
12:AL:82:VAL:HG22	12:AL:83:LEU:H	1.83	0.42
30:BH:88:ILE:HG13	30:BH:144:VAL:HG11	2.02	0.42
43:BU:31:LEU:N	43:BU:31:LEU:HD23	2.26	0.42
23:DA:1541:U:C3'	23:DA:1542:G:O3'	2.65	0.42
40:DR:77:ALA:C	40:DR:79:VAL:N	2.73	0.42
24:BB:73:A:H3'	24:BB:74:U:C6	2.54	0.42
23:BA:2777:G:H8	23:BA:2777:G:H3'	1.84	0.42
23:DA:2777:G:H8	23:DA:2777:G:H3'	1.85	0.42
46:BX:35:THR:HB	46:BX:36:GLY:H	1.54	0.42
23:DA:102:G:C8	23:DA:102:G:H5''	2.55	0.42
1:AA:1149:C:O5'	1:AA:1149:C:H6	2.02	0.42
28:DF:129:GLY:HA3	28:DF:163:ALA:HB3	2.01	0.42
23:BA:301:G:H5'	23:BA:334:C:O2'	2.19	0.42
23:DA:1494:A:HO2'	23:DA:1495:A:P	2.42	0.42
1:CA:554:C:H2'	1:CA:555:C:C6	2.53	0.42
35:BM:66:ILE:HG22	35:BM:104:PHE:HD2	1.82	0.42
44:DV:6:LYS:HG3	44:DV:8:TYR:CZ	2.54	0.42
23:DA:55:G:H2'	23:DA:56:A:C8	2.54	0.42
1:CA:1004:A:H3'	1:CA:1004:A:N3	2.35	0.42
23:BA:1833:U:N3	23:BA:1834:U:C5	2.88	0.42
23:BA:2572:A:C8	26:BD:144:ARG:HB3	2.54	0.42
1:CA:502:G:H2'	1:CA:503:C:O4'	2.20	0.42
23:BA:150:C:H2'	23:BA:151:C:H6	1.83	0.42
23:BA:1324:G:H4'	23:BA:1616:A:C2	2.54	0.42
53:D5:29:LYS:O	53:D5:29:LYS:HG2	2.19	0.42
23:BA:909:A:C4	23:BA:912:C:C5	3.07	0.42
36:BN:100:LEU:HD21	36:BN:113:LEU:HB2	2.02	0.42
6:AF:82:ARG:HB2	6:AF:85:VAL:HG23	2.01	0.42
23:DA:2820:A:O4'	36:DN:5:LYS:HG3	2.18	0.42
2:AB:32:ILE:HA	2:AB:32:ILE:HD12	1.89	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:CG:95:ARG:O	7:CG:99:LEU:HG	2.20	0.42
1:CA:619:U:C2	4:CD:135:LEU:HD21	2.54	0.42
23:DA:1414:G:C4	23:DA:1415:U:C5	3.07	0.42
23:DA:2096:U:H2'	23:DA:2097:C:C6	2.55	0.42
23:BA:1648:C:H2'	23:BA:1649:G:O5'	2.19	0.42
23:BA:2716:U:O2'	23:BA:2717:G:H5'	2.19	0.42
1:AA:1020:U:H2'	1:AA:1021:G:H8	1.85	0.42
11:AK:48:ILE:HD11	11:AK:64:ALA:HA	2.00	0.42
28:BF:34:LEU:HD21	28:BF:159:VAL:HG21	2.00	0.42
7:CG:87:VAL:HG11	7:CG:154:TYR:O	2.20	0.42
23:BA:1109:C:H2'	23:BA:1110:G:O4'	2.19	0.42
23:BA:1229:G:H2'	23:BA:1230:C:C6	2.54	0.42
1:CA:541:G:O2'	4:CD:41:GLY:HA2	2.19	0.42
23:DA:1285:G:C5	23:DA:1329:U:C4	3.07	0.42
23:BA:1954:G:N2	23:BA:1956:U:H3	2.18	0.42
1:AA:1323:G:H2'	1:AA:1324:A:C8	2.55	0.42
23:DA:2212:A:HO2'	23:DA:2215:G:C1'	2.32	0.42
1:AA:1145:C:H4'	1:AA:1146:A:H8	1.85	0.42
8:CH:84:ARG:O	8:CH:135:CYS:HB2	2.19	0.42
26:DD:17:ASP:OD1	26:DD:18:ASP:N	2.53	0.42
50:D2:9:LYS:HA	50:D2:9:LYS:HD3	1.88	0.42
26:DD:197:ILE:HD11	26:DD:199:ARG:HE	1.84	0.42
28:DF:178:PHE:HA	28:DF:179:PRO:HD3	1.79	0.42
23:DA:1399:C:O2'	23:DA:1400:G:H5'	2.17	0.42
5:CE:147:ASP:O	5:CE:151:LEU:HG	2.20	0.42
44:DV:48:PHE:CE2	44:DV:71:VAL:HG21	2.55	0.42
23:BA:846:C:C2	23:BA:847:U:C5	3.07	0.42
23:DA:947:G:H2'	23:DA:948:G:C5'	2.43	0.42
1:AA:1378:C:H5	1:AA:1379:G:C8	2.36	0.42
23:DA:2516:G:C6	23:DA:2517:C:C4	3.07	0.42
30:DH:82:ARG:HB3	30:DH:89:TYR:CG	2.55	0.42
23:BA:643:A:O2'	23:BA:644:A:H5'	2.19	0.42
23:BA:1541:U:C3'	23:BA:1542:G:O3'	2.66	0.42
23:DA:646:A:H2'	23:DA:647:G:O4'	2.20	0.42
27:BE:160:ASN:ND2	27:BE:162:LEU:H	2.18	0.42
1:AA:769:G:H4'	1:AA:1513:A:H4'	2.02	0.42
26:DD:4:ILE:HG13	26:DD:28:ALA:HB1	2.01	0.42
23:DA:2850:A:H2'	23:DA:2851:A:O4'	2.19	0.42
23:DA:1265:A:H3'	50:D2:19:ARG:NH1	2.35	0.42
1:CA:1149:C:O5'	1:CA:1149:C:H6	2.02	0.42
23:BA:332:A:O2'	23:BA:334:C:OP2	2.30	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DA:1495:A:N3	23:DA:1496:A:C2	2.87	0.42
44:BV:9:TYR:O	44:BV:38:TYR:HE2	2.02	0.42
41:BS:4:LYS:HA	41:BS:106:ILE:HG22	2.02	0.42
35:BM:26:TYR:O	35:BM:26:TYR:CD1	2.72	0.42
1:CA:1004:A:H8	1:CA:1026:G:N7	2.17	0.42
3:CC:173:VAL:O	3:CC:173:VAL:HG12	2.19	0.42
23:DA:2196:C:O2'	23:DA:2197:U:H5'	2.19	0.42
35:DM:68:ILE:HD13	35:DM:103:MET:HG3	2.02	0.42
23:DA:1324:G:H4'	23:DA:1616:A:C2	2.54	0.42
26:BD:11:MET:CB	26:BD:24:THR:HA	2.50	0.42
2:CB:32:ILE:HD12	2:CB:32:ILE:HA	1.90	0.42
26:DD:11:MET:CB	26:DD:24:THR:HA	2.49	0.42
1:AA:284:G:H2'	1:AA:285:G:C8	2.54	0.42
1:CA:1323:G:H2'	1:CA:1324:A:C8	2.55	0.42
1:AA:1234:C:O2'	1:AA:1235:U:H5'	2.20	0.42
7:AG:87:VAL:HA	7:AG:88:PRO:HD3	1.80	0.42
24:BB:16:G:C6	24:BB:69:G:C2	3.08	0.42
4:AD:78:LEU:O	4:AD:81:GLU:HB3	2.18	0.42
23:DA:1759:A:C8	23:DA:2696:U:H1'	2.54	0.42
13:AM:24:GLY:O	13:AM:25:ILE:HD13	2.19	0.42
2:CB:15:VAL:C	2:CB:16:HIS:CG	2.93	0.42
23:BA:1759:A:C8	23:BA:2696:U:H1'	2.53	0.42
23:DA:1546:A:H2'	23:DA:154(B):C:O4'	2.20	0.42
1:AA:1259:C:O5'	1:AA:1259:C:H6	2.03	0.42
27:BE:9:ILE:H	27:BE:9:ILE:HD13	1.83	0.42
10:AJ:79:ARG:HD3	10:AJ:79:ARG:HA	1.88	0.42
26:BD:110:GLY:HA2	26:BD:162:ALA:HB2	2.01	0.42
23:BA:335:C:C2	23:BA:336:C:C5	3.07	0.42
39:BQ:79:PHE:HE1	39:BQ:83:LEU:CD2	2.32	0.42
40:BR:40:LEU:C	40:BR:45:THR:HB	2.40	0.42
40:BR:45:THR:O	40:BR:46:VAL:C	2.56	0.42
23:DA:1828:G:OP2	25:DC:239:ARG:NH2	2.53	0.42
35:DM:75:THR:HG21	35:DM:85:LYS:HZ2	1.82	0.42
42:DT:30:VAL:HG12	42:DT:31:HIS:H	1.84	0.42
1:AA:980:C:O2	14:AN:19:ARG:HA	2.18	0.42
49:B1:42:CYS:HA	49:B1:59:VAL:C	2.40	0.42
35:DM:141:GLN:NE2	44:DV:72:ARG:HG2	2.35	0.42
35:BM:141:GLN:OXT	44:BV:98:MET:HE3	2.19	0.42
5:AE:33:VAL:CG1	5:AE:109:ILE:HD13	2.50	0.42
23:DA:1310:G:C3'	23:DA:1311:G:H5''	2.49	0.42
23:DA:1310:G:H2'	23:DA:1311:G:H5''	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:DD:101:ARG:HB3	26:DD:169:ASN:HD22	1.83	0.42
3:CC:23:TYR:HB2	10:CJ:93:GLY:O	2.19	0.42
23:DA:778:G:C5	23:DA:779:U:C4	3.07	0.42
29:BG:46:GLU:HG3	29:BG:51:ARG:CD	2.49	0.42
7:AG:15:ASP:HB3	7:AG:19:GLY:N	2.28	0.42
23:DA:94:G:N3	47:DY:47:ASN:ND2	2.67	0.42
34:DL:9:ASN:N	34:DL:10:PRO:CD	2.81	0.42
2:CB:11:LEU:HD12	2:CB:217:ARG:NH2	2.34	0.42
23:DA:85:G:C8	23:DA:85:G:H5''	2.50	0.42
23:BA:2809:A:N6	23:BA:2892:A:C8	2.88	0.42
23:DA:840:C:H2'	23:DA:841:A:C8	2.54	0.42
28:DF:115:ARG:CD	28:DF:115:ARG:H	2.30	0.42
1:CA:222:U:H2'	1:CA:223:U:H6	1.83	0.42
1:CA:1116:C:C3'	1:CA:1117:G:H5''	2.49	0.42
25:DC:126:GLN:HG2	25:DC:127:VAL:H	1.85	0.42
23:BA:1996:C:H5	33:BK:32:TYR:OH	2.03	0.42
23:DA:2749:A:H1'	29:DG:63:SER:OG	2.20	0.42
1:AA:192:U:H2'	1:AA:193:C:C6	2.55	0.42
23:DA:297:C:H2'	23:DA:298:G:O4'	2.20	0.42
35:BM:74:TYR:O	35:BM:89:ASN:N	2.53	0.42
1:CA:1201:A:O2'	1:CA:1202:G:OP2	2.34	0.42
30:BH:118:LYS:HA	30:BH:119:PRO:HD3	1.86	0.42
10:AJ:53:PRO:HA	14:AN:42:ILE:HD11	2.02	0.42
26:DD:107:THR:O	26:DD:190:GLY:HA2	2.20	0.42
1:AA:338:A:C6	1:AA:339:C:N4	2.87	0.42
23:BA:2767:C:H2'	23:BA:2768:C:H6	1.84	0.42
15:CO:8:LYS:O	15:CO:12:ILE:HG13	2.19	0.42
23:BA:536:A:H2'	23:BA:537:C:C6	2.55	0.42
23:DA:648:G:O4'	23:DA:2351:G:H5''	2.20	0.42
23:DA:2870:C:H2'	23:DA:2871:C:O4'	2.20	0.42
23:BA:540:G:C4	23:BA:541:C:C5	3.08	0.42
35:DM:62:GLY:HA2	44:DV:116:VAL:HG22	2.02	0.42
23:BA:2212:A:HO2'	23:BA:2215:G:C1'	2.32	0.42
35:BM:18:LYS:HB3	35:BM:19:GLY:H	1.63	0.42
1:CA:1466:C:H2'	1:CA:1467:G:O4'	2.20	0.42
23:DA:1918:A:O2'	23:DA:1920:C:N4	2.53	0.42
13:CM:24:GLY:O	13:CM:25:ILE:HD13	2.20	0.42
1:CA:666:G:C5	1:CA:741:G:C6	3.08	0.42
14:CN:12:ARG:HB3	14:CN:14:PRO:HD3	2.02	0.42
4:AD:199:ASN:ND2	4:AD:202:LEU:HG	2.35	0.42
25:BC:218:ARG:HH11	25:BC:218:ARG:HG3	1.84	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
44:BV:70:LEU:HD23	44:BV:70:LEU:N	2.34	0.42
11:AK:38:ASN:ND2	11:AK:38:ASN:N	2.67	0.42
33:DK:19:ILE:HD13	33:DK:19:ILE:H	1.84	0.42
23:DA:2322:A:H3'	23:DA:2323:G:H8	1.83	0.42
6:AF:3:ARG:HD3	6:AF:64:GLN:OE1	2.20	0.42
23:DA:2618:G:C6	23:DA:2619:C:C4	3.07	0.42
23:DA:2070:G:H2'	23:DA:2071:A:O4'	2.19	0.42
1:AA:1399:C:H4'	1:AA:1400:C:C5'	2.20	0.42
23:BA:588:U:H2'	23:BA:589:C:H6	1.82	0.42
34:BL:47:ASP:HB3	34:BL:48:PRO:CA	2.49	0.42
23:BA:1828:G:OP2	25:BC:239:ARG:NH2	2.53	0.42
34:DL:86:LYS:HG3	34:DL:87:ASP:N	2.34	0.42
23:BA:72:U:O2'	23:BA:73:A:H5'	2.19	0.42
23:BA:2711:A:OP1	23:BA:712(B):A:P	2.78	0.42
25:DC:32:SER:HA	25:DC:36:PRO:HG2	2.01	0.42
10:AJ:32:ALA:HB2	10:AJ:76:ASN:HB2	2.00	0.42
26:BD:91:VAL:CB	26:BD:95:ILE:HD11	2.43	0.42
23:BA:2262:U:H4'	23:BA:2328:A:C2	2.55	0.42
28:DF:111:LEU:N	28:DF:112:PRO:CD	2.83	0.42
30:DH:133:HIS:HA	30:DH:134:PRO:HD3	1.90	0.42
25:BC:31:LYS:HG3	25:BC:33:LEU:HG	2.02	0.42
25:BC:27:THR:CG2	25:BC:83:GLU:HG2	2.50	0.42
12:AL:74:HIS:HD2	12:AL:76:LEU:HB2	1.84	0.42
1:CA:1328:C:H5''	13:CM:28:ALA:HB1	2.02	0.42
16:CP:20:VAL:HG22	16:CP:21:VAL:N	2.34	0.42
30:DH:88:ILE:HG13	30:DH:144:VAL:HG11	2.02	0.42
34:DL:30:THR:CG2	34:DL:31:ALA:N	2.80	0.42
1:CA:392:G:C2	1:CA:393:A:C5	3.08	0.42
25:BC:147:LEU:HD13	25:BC:155:LEU:CD1	2.45	0.42
4:CD:51:PRO:HB3	4:CD:55:ALA:CB	2.49	0.42
46:BX:43:TYR:HA	46:BX:44:PRO:HD3	1.77	0.42
23:BA:2484:G:C2	23:BA:2485:G:C8	3.07	0.42
23:BA:804:A:H5''	23:BA:805:G:OP1	2.20	0.42
23:BA:1678:G:O5'	23:BA:1678:G:C8	2.64	0.42
39:BQ:62:ILE:HD12	39:BQ:76:TYR:CZ	2.54	0.42
51:B3:34:LEU:H	51:B3:34:LEU:HD13	1.85	0.42
25:BC:79:VAL:HG11	25:BC:111:LEU:CD1	2.50	0.42
36:DN:100:LEU:HD21	36:DN:113:LEU:HB2	2.01	0.42
34:BL:7:ARG:O	34:BL:10:PRO:HD3	2.19	0.42
1:CA:1528:U:H6	1:CA:1528:U:C5'	2.31	0.42
23:DA:839:U:H2'	23:DA:840:C:H6	1.82	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BA:426:C:C2	23:BA:427:U:C5	3.08	0.42
47:BY:56:GLN:O	47:BY:60:LEU:HG	2.20	0.42
1:AA:222:U:H2'	1:AA:223:U:H6	1.83	0.42
41:DS:4:LYS:HA	41:DS:106:ILE:HG22	2.01	0.42
1:CA:192:U:H2'	1:CA:193:C:C6	2.55	0.42
6:CF:8:ILE:HD11	6:CF:79:LEU:HD13	2.01	0.42
23:DA:919:G:H2'	23:DA:920:G:H8	1.85	0.42
23:BA:919:G:H2'	23:BA:920:G:H8	1.85	0.42
1:CA:1281:U:O2'	1:CA:1282:C:P	2.78	0.42
23:DA:628:G:C6	23:DA:636:G:C2	3.07	0.42
23:BA:56:A:C2	23:BA:115:C:O2	2.73	0.42
3:AC:33:LEU:HD21	14:AN:53:LEU:HD23	2.00	0.42
25:DC:166:GLN:HB2	25:DC:174:ILE:HG22	2.00	0.42
23:BA:2090:G:H21	46:BX:45:ASN:HD21	1.67	0.42
23:BA:363(D):G:H2'	23:BA:363(E):G:C8	2.55	0.42
23:DA:1040:C:H2'	23:DA:1041:C:H6	1.85	0.42
23:DA:433:C:O2'	23:DA:434:U:H5'	2.19	0.42
1:AA:338:A:C6	1:AA:339:C:C4	3.08	0.42
41:BS:107:LEU:CD1	41:BS:107:LEU:N	2.82	0.42
29:BG:96:ALA:CB	29:BG:105:LEU:HB3	2.49	0.42
1:CA:338:A:C6	1:CA:339:C:C4	3.08	0.42
23:BA:286:C:C2	23:BA:287:C:C5	3.07	0.42
35:BM:130:LYS:HG2	35:BM:131:ILE:N	2.33	0.42
23:DA:1695:G:N2	23:DA:1696:G:C8	2.88	0.42
20:CT:94:ALA:C	20:CT:96:GLY:H	2.23	0.42
6:CF:3:ARG:HD3	6:CF:64:GLN:OE1	2.20	0.42
32:BJ:35:ARG:O	32:BJ:73:ASP:HB3	2.20	0.42
7:CG:22:LEU:HD23	7:CG:63:LYS:HZ2	1.85	0.42
23:BA:665:C:O2'	23:BA:666:G:H5'	2.20	0.42
41:DS:30:GLU:HA	41:DS:33:ARG:HD2	2.01	0.42
39:DQ:57:PHE:HB3	39:DQ:61:TRP:CZ2	2.55	0.42
23:DA:191:A:H2'	23:DA:192:C:C6	2.55	0.42
1:AA:1419:G:C6	1:AA:1420:C:C4	3.07	0.42
23:DA:460:A:C6	23:DA:470:A:C8	3.08	0.42
7:AG:23:VAL:HG13	7:AG:43:PHE:CE2	2.55	0.42
26:BD:97:LYS:HE2	26:BD:97:LYS:HB3	1.91	0.42
50:D2:13:LYS:HB3	50:D2:13:LYS:HE2	1.85	0.42
24:DB:3:C:H2'	24:DB:4:C:C6	2.55	0.42
1:CA:1419:G:C6	1:CA:1420:C:C4	3.08	0.42
2:AB:15:VAL:C	2:AB:16:HIS:CG	2.92	0.42
9:AI:33:PHE:HZ	9:AI:43:ALA:O	2.03	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BA:2444:G:C6	23:BA:2445:G:C5	3.08	0.42
23:DA:2444:G:C6	23:DA:2445:G:C5	3.07	0.42
34:DL:51:PHE:HB3	34:DL:52:GLU:H	1.76	0.42
25:DC:222:ARG:HH12	25:DC:239:ARG:CZ	2.33	0.42
32:BJ:40:ASP:OD1	32:BJ:42:GLU:HG2	2.20	0.42
23:DA:114(B):A:C2	23:DA:1144:G:C8	3.08	0.42
47:DY:2:LYS:HD3	47:DY:2:LYS:H	1.84	0.42
35:DM:81:VAL:HG12	35:DM:82:ARG:CG	2.41	0.42
1:CA:1346:A:H5'	9:CI:120:ARG:NH1	2.25	0.42
23:BA:948:G:OP1	23:BA:962:G:OP1	2.38	0.42
28:BF:111:LEU:N	28:BF:112:PRO:CD	2.83	0.42
3:CC:182:ILE:HG23	3:CC:202:ILE:C	2.40	0.42
23:DA:573:G:O2'	23:DA:574:C:H3'	2.20	0.42
43:DU:27:VAL:HG23	43:DU:27:VAL:O	2.19	0.42
26:DD:171:GLU:HG2	26:DD:185:LYS:CG	2.50	0.42
2:CB:187:LEU:CD1	2:CB:205:ASP:HB3	2.50	0.42
3:CC:59:ARG:HA	3:CC:63:ASN:O	2.20	0.42
23:DA:2468:G:HO2'	23:DA:2469:A:P	2.41	0.42
30:BH:101:LEU:HG	30:BH:107:ILE:CG2	2.49	0.42
4:CD:49:ARG:HA	4:CD:49:ARG:HD2	1.89	0.42
23:DA:479:A:H4'	23:DA:480:A:O5'	2.20	0.42
23:DA:1349:A:N6	23:DA:1598:C:H42	2.17	0.42
1:AA:586:C:O2'	1:AA:878:G:H4'	2.20	0.42
41:BS:18:ARG:HG2	41:BS:76:VAL:HG11	2.02	0.42
36:BN:2:ARG:HB3	36:BN:3:HIS:CE1	2.54	0.42
27:BE:106:ARG:HG2	27:BE:106:ARG:H	1.54	0.42
5:AE:92:LYS:O	5:AE:118:ILE:HD12	2.20	0.42
40:BR:28:GLU:OE1	40:BR:31:ALA:HB2	2.19	0.42
23:DA:2477:C:HO2'	23:DA:2478:A:P	2.43	0.42
1:CA:235:C:H1'	17:CQ:61:GLU:CD	2.40	0.42
1:AA:663:A:O2'	1:AA:664:G:H5'	2.20	0.42
23:BA:514:A:H1'	23:BA:581:C:O2'	2.19	0.42
6:AF:8:ILE:HD11	6:AF:79:LEU:HD13	2.02	0.42
1:AA:815:A:H4'	1:AA:817:C:C5	2.55	0.42
3:AC:173:VAL:HG12	3:AC:173:VAL:O	2.20	0.42
23:DA:2341:G:H2'	23:DA:2342:C:C6	2.54	0.42
23:BA:2536:G:C6	23:BA:2537:U:C4	3.08	0.42
25:BC:105:ILE:HG13	25:BC:106:ILE:HD12	2.02	0.42
25:BC:260:ARG:O	25:BC:261:LYS:C	2.58	0.42
45:DW:82:ARG:HA	45:DW:83:PRO:HD2	1.91	0.42
23:BA:1751:C:H2'	23:BA:1752:C:C6	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:DC:261:LYS:HB2	25:DC:261:LYS:NZ	2.35	0.42
23:BA:1465:G:C2	23:BA:1466:G:C8	3.07	0.42
1:CA:156:G:C2	1:CA:166:G:C2	3.07	0.42
52:D4:47:ARG:O	52:D4:48:LYS:HB2	2.19	0.42
23:DA:335:C:C2	23:DA:336:C:C5	3.08	0.42
26:BD:17:ASP:OD1	26:BD:18:ASP:N	2.52	0.42
38:BP:131:ALA:O	38:BP:135:VAL:HG23	2.20	0.42
23:DA:173:G:H2'	23:DA:174:C:C6	2.54	0.42
23:BA:1107:G:O2'	23:BA:1108:U:H5'	2.19	0.42
23:DA:924:C:H2'	23:DA:925:C:C6	2.54	0.42
43:BU:34:LYS:HE2	43:BU:34:LYS:HB3	1.80	0.42
29:DG:144:VAL:O	29:DG:148:ILE:HG12	2.20	0.42
32:DJ:35:ARG:O	32:DJ:73:ASP:HB3	2.20	0.42
39:DQ:92:ARG:CD	39:DQ:95:LEU:H	2.32	0.42
39:DQ:92:ARG:HH21	40:DR:11:GLN:H	1.67	0.42
39:BQ:83:LEU:HG	39:BQ:88:ILE:CD1	2.43	0.42
1:CA:1130:A:H4'	9:CI:20:ARG:HH22	1.84	0.42
15:AO:36:ILE:HD12	15:AO:63:ARG:NH1	2.33	0.42
5:AE:76:ILE:CG1	5:AE:77:PRO:HD2	2.31	0.42
42:BT:26:TYR:CE1	42:BT:83:VAL:HG21	2.55	0.42
42:DT:26:TYR:CE1	42:DT:83:VAL:HG21	2.55	0.42
23:DA:2247:A:H2'	23:DA:2248:C:H6	1.83	0.42
53:B5:25:MET:SD	53:B5:47:LYS:HG2	2.59	0.42
53:B5:33:ASN:ND2	53:B5:34:TRP:N	2.67	0.42
5:AE:147:ASP:O	5:AE:151:LEU:HG	2.20	0.42
43:BU:30:VAL:HG13	43:BU:37:VAL:HG12	2.01	0.42
44:BV:97:GLU:HB3	44:BV:125:LEU:CD2	2.43	0.42
23:DA:2262:U:H4'	23:DA:2328:A:C2	2.55	0.42
25:DC:35:LYS:HG3	25:DC:104:TYR:CD2	2.55	0.42
25:DC:35:LYS:CB	25:DC:36:PRO:HD3	2.50	0.42
23:BA:1579:A:H2'	23:BA:1580:A:C8	2.55	0.42
12:CL:69:ILE:HA	12:CL:99:ILE:HG22	2.01	0.42
23:BA:307:G:N1	23:BA:310:A:OP2	2.53	0.42
4:CD:31:CYS:O	4:CD:32:ALA:CB	2.68	0.42
25:DC:76:PRO:HA	25:DC:118:VAL:HG23	2.02	0.42
23:BA:464:U:H4'	52:B4:5:TRP:CZ3	2.54	0.42
44:BV:179:ASP:CG	44:BV:180:VAL:N	2.73	0.42
4:AD:122:ARG:O	4:AD:134:ASP:HB2	2.19	0.42
1:AA:1428:A:H2'	1:AA:1429:C:C6	2.55	0.42
23:BA:1349:A:N6	23:BA:1598:C:H42	2.18	0.42
4:CD:104:VAL:HG11	4:CD:146:ILE:HD12	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:AI:14:VAL:HG12	9:AI:15:ALA:N	2.35	0.42
2:AB:71:VAL:HG23	2:AB:164:VAL:HG13	2.02	0.42
23:BA:1951:U:O2	23:BA:1953:A:H8	2.03	0.42
23:BA:2686:G:C5	23:BA:2687:U:C4	3.08	0.42
28:BF:128:ARG:HH21	28:BF:130:ASN:ND2	2.17	0.42
23:DA:2809:A:N6	23:DA:2892:A:C8	2.88	0.42
1:CA:948:C:OP1	13:CM:107:ALA:HA	2.20	0.42
46:DX:37:ILE:HG23	46:DX:38:SER:N	2.35	0.42
25:BC:145:VAL:HG12	25:BC:146:GLU:O	2.20	0.42
1:AA:502:G:OP1	12:AL:115:SER:HA	2.19	0.42
47:DY:21:LEU:CD2	47:DY:22:GLU:HG3	2.50	0.42
1:AA:804:U:H5''	1:AA:805:C:OP2	2.20	0.42
23:BA:1655:A:O2'	26:BD:115:GLY:HA2	2.19	0.42
23:DA:1862:G:H2'	23:DA:1863:G:C8	2.53	0.42
35:DM:74:TYR:O	35:DM:89:ASN:N	2.52	0.42
51:B3:30:THR:CG2	51:B3:31:PRO:HD2	2.48	0.42
19:CS:53:ASN:ND2	19:CS:56:GLN:H	2.17	0.42
25:BC:261:LYS:NZ	25:BC:261:LYS:HB2	2.35	0.42
25:DC:105:ILE:HG12	25:DC:106:ILE:HD12	2.02	0.42
18:CR:45:SER:HB3	18:CR:51:LEU:CG	2.49	0.42
1:AA:1330:U:H4'	13:AM:23:TYR:CE2	2.54	0.42
41:BS:70:TYR:HD2	41:BS:70:TYR:H	1.68	0.42
1:CA:862:C:H2'	1:CA:863:U:O4'	2.20	0.42
23:BA:1464:C:O2	23:BA:1528:A:H2	2.03	0.42
23:BA:173:G:H2'	23:BA:174:C:C6	2.55	0.42
15:CO:7:GLU:HA	15:CO:10:LYS:HB3	2.01	0.42
33:BK:17:ARG:HB2	33:BK:45:GLU:HG3	2.02	0.42
23:BA:2396:G:O2'	23:BA:2397:G:H5'	2.20	0.42
38:BP:23:ARG:NH2	38:BP:120:ARG:HD3	2.35	0.42
14:AN:12:ARG:HB3	14:AN:14:PRO:HD3	2.02	0.42
1:AA:69:G:H2'	1:AA:73:G:H8	1.84	0.42
33:DK:17:ARG:HB2	33:DK:45:GLU:HG3	2.01	0.42
23:DA:1726:G:C2	23:DA:1735:U:O2	2.73	0.42
39:DQ:59:ARG:O	39:DQ:63:VAL:HG23	2.20	0.42
7:AG:25:ALA:O	7:AG:29:LYS:HG2	2.20	0.42
29:BG:12:PRO:O	29:BG:15:VAL:HG22	2.20	0.42
1:CA:1020:U:H2'	1:CA:1021:G:H8	1.85	0.42
1:CA:119:A:C5	1:CA:240:C:C4	3.07	0.42
2:CB:113:HIS:O	2:CB:116:GLU:HG2	2.19	0.42
44:DV:157:LEU:N	44:DV:157:LEU:HD12	2.34	0.42
23:BA:2194:G:H2'	23:BA:2195:C:H6	1.84	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:272:C:H2'	1:AA:273:A:H8	1.85	0.42
7:CG:111:ARG:HA	7:CG:112:PRO:HD3	1.82	0.42
40:BR:4:ILE:HG22	40:BR:5:VAL:N	2.35	0.42
23:DA:1140:C:OP1	32:DJ:46:LEU:HB3	2.20	0.42
53:D5:33:ASN:ND2	53:D5:34:TRP:N	2.67	0.42
49:D1:50:THR:CG2	49:D1:51:TYR:H	2.14	0.42
34:BL:85:LEU:HD21	34:BL:116:GLY:O	2.20	0.42
42:DT:31:HIS:HA	42:DT:32:PRO:HD3	1.97	0.42
37:DO:49:VAL:HG13	37:DO:76:LYS:HB2	2.02	0.42
25:BC:190:TYR:O	25:BC:191:ALA:HB2	2.20	0.42
25:DC:197:GLY:O	25:DC:198:ASN:C	2.58	0.42
10:CJ:33:GLN:HB2	10:CJ:75:ILE:CD1	2.50	0.42
25:DC:94:LEU:HB2	25:DC:104:TYR:CE1	2.47	0.42
23:DA:2028:U:O4	23:DA:2033:A:OP1	2.38	0.42
34:BL:61:ARG:HH11	53:B5:13:ARG:HD2	1.84	0.42
7:AG:70:LYS:HA	7:AG:71:PRO:HD2	1.94	0.42
1:CA:1328:C:H5'	13:CM:28:ALA:CB	2.49	0.42
4:CD:50:ARG:HA	4:CD:51:PRO:HD3	1.85	0.42
23:BA:1334:G:C6	23:BA:1335:U:C4	3.07	0.42
28:BF:18:GLU:HB3	28:BF:175:LEU:HD13	2.02	0.42
28:DF:161:THR:HG21	28:DF:172:LEU:CD2	2.50	0.42
1:AA:391:G:C6	1:AA:392:G:C5	3.08	0.42
1:AA:429:U:H4'	1:AA:430:A:O5'	2.20	0.42
23:BA:1447:G:N3	23:BA:1545:A:H2	2.18	0.42
23:BA:388:G:OP1	46:BX:33:LYS:CB	2.68	0.42
23:DA:907:U:O2'	35:DM:101:ARG:NH2	2.49	0.42
3:CC:34:LEU:HD21	3:CC:38:ARG:HH21	1.84	0.42
24:BB:106:G:C5	24:BB:107:U:C5	3.08	0.42
1:AA:1513:A:H2'	1:AA:1514:C:C6	2.55	0.42
4:CD:108:LEU:HB3	4:CD:110:PHE:HD2	1.85	0.42
23:DA:595:C:H2'	23:DA:596:G:O4'	2.20	0.42
28:BF:128:ARG:HG2	28:BF:129:GLY:H	1.85	0.42
23:DA:1602:U:H3'	23:DA:1603:A:C5'	2.50	0.42
11:AK:59:TYR:O	11:AK:62:GLN:HB3	2.19	0.42
1:CA:685:G:O2'	1:CA:686:U:H5'	2.20	0.42
23:BA:311:A:C8	23:BA:332:A:C5	3.07	0.42
33:BK:22:ILE:HG12	33:BK:42:SER:H	1.85	0.42
1:AA:805:C:H2'	1:AA:806:C:C6	2.53	0.42
1:AA:1004:A:H3'	1:AA:1004:A:N3	2.34	0.42
23:DA:56:A:C2	23:DA:115:C:O2	2.73	0.42
6:CF:75:LEU:C	6:CF:75:LEU:HD23	2.40	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:BD:144:ARG:HB3	26:BD:145:LYS:H	1.57	0.42
23:BA:2050:C:H1'	26:BD:156:MET:CE	2.49	0.42
20:CT:63:ILE:HG21	20:CT:81:LYS:HG3	2.00	0.42
10:AJ:30:SER:HB2	10:AJ:80:LYS:HG3	2.02	0.42
30:DH:29:TYR:C	30:DH:32:PRO:HD2	2.40	0.42
1:CA:1432:G:OP1	38:DP:107:ASP:HB2	2.20	0.42
18:AR:45:SER:HB3	18:AR:51:LEU:HG	2.02	0.42
19:CS:51:VAL:HG12	19:CS:52:TYR:O	2.20	0.42
23:BA:161:U:HO2'	23:BA:162:U:H5	1.66	0.42
23:DA:912:C:C2	23:DA:913:U:C5	3.08	0.42
6:CF:82:ARG:HB2	6:CF:85:VAL:HG23	2.01	0.42
23:DA:2886:G:H2'	23:DA:2887:U:H6	1.84	0.42
23:BA:2590:A:C2	23:BA:2605:U:C2	3.08	0.42
41:DS:107:LEU:CD1	41:DS:107:LEU:N	2.83	0.42
5:AE:35:GLY:HA3	5:AE:41:VAL:HG12	2.02	0.42
1:AA:173:U:C6	1:AA:197:A:C2	3.08	0.42
1:CA:1020:U:H2'	1:CA:1021:G:C8	2.55	0.42
41:BS:30:GLU:HA	41:BS:33:ARG:HD2	2.02	0.42
23:BA:2762:G:O2'	23:BA:2763:G:H5'	2.20	0.42
23:DA:699:A:H4'	23:DA:1634:A:N7	2.34	0.42
40:BR:1:MET:SD	40:BR:42:GLY:HA3	2.60	0.42
24:BB:33:G:O2'	24:BB:34:U:H5'	2.20	0.42
23:BA:924:C:H2'	23:BA:925:C:C6	2.55	0.42
23:BA:2870:C:H2'	23:BA:2871:C:O4'	2.19	0.42
1:CA:386:C:H2'	1:CA:387:U:H5''	2.02	0.42
23:DA:284:U:H2'	23:DA:285:C:C6	2.55	0.42
8:CH:85:ARG:HH12	8:CH:134:ILE:HG23	1.85	0.42
28:DF:34:LEU:HD21	28:DF:159:VAL:HG21	2.02	0.42
38:DP:131:ALA:O	38:DP:135:VAL:HG23	2.20	0.42
52:D4:18:PHE:CD2	52:D4:18:PHE:C	2.93	0.42
7:AG:60:LYS:HA	7:AG:60:LYS:HD2	1.91	0.42
23:DA:2284:C:O5'	23:DA:2284:C:H6	2.03	0.42
29:DG:78:GLY:HA2	29:DG:83:TYR:CE1	2.54	0.42
44:DV:40:ASP:OD1	44:DV:42:VAL:HG12	2.19	0.42
2:AB:69:LEU:HD22	2:AB:159:PRO:HG2	2.02	0.41
34:BL:62:LEU:HD22	34:BL:62:LEU:N	2.35	0.41
5:CE:20:GLN:O	5:CE:21:ALA:C	2.59	0.41
37:DO:28:VAL:HG13	37:DO:35:ILE:HD11	2.02	0.41
49:D1:42:CYS:HA	49:D1:59:VAL:C	2.41	0.41
23:DA:270(K):G:H2'	23:DA:270(L):C:O4'	2.20	0.41
25:DC:35:LYS:HZ1	25:DC:104:TYR:H	1.68	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:CC:137:ALA:O	3:CC:141:VAL:HG23	2.20	0.41
2:AB:187:LEU:CD1	2:AB:205:ASP:HB3	2.50	0.41
2:AB:80:ILE:HD12	2:AB:211:ILE:HB	2.01	0.41
7:AG:92:SER:O	7:AG:96:GLN:HG3	2.20	0.41
12:CL:74:HIS:CD2	12:CL:76:LEU:HB2	2.55	0.41
12:CL:82:VAL:HG23	12:CL:106:ALA:CB	2.45	0.41
26:DD:116:VAL:HG13	26:DD:117:MET:N	2.34	0.41
38:DP:61:PHE:CZ	38:DP:76:PHE:HB2	2.54	0.41
24:BB:72:G:N2	24:BB:103:U:C5	2.88	0.41
19:CS:29:ARG:HB2	19:CS:48:THR:N	2.32	0.41
28:BF:15:VAL:HG22	28:BF:175:LEU:HB3	2.02	0.41
23:DA:1543:A:H3'	23:DA:1543:A:C8	2.55	0.41
23:DA:65:C:O2'	23:DA:66:C:H5'	2.20	0.41
47:BY:46:GLN:O	47:BY:47:ASN:CB	2.61	0.41
25:DC:235:GLY:O	25:DC:237:GLU:N	2.48	0.41
23:BA:811:U:H3'	34:BL:25:SER:O	2.20	0.41
7:CG:150:ALA:HA	11:CK:59:TYR:HD2	1.85	0.41
23:BA:2809:A:N1	23:BA:2892:A:C4	2.88	0.41
23:DA:363(A):G:H2'	23:DA:363(B):A:C8	2.54	0.41
13:CM:106:ASN:HB2	13:CM:107:ALA:H	1.55	0.41
7:AG:150:ALA:HA	11:AK:59:TYR:CD2	2.55	0.41
38:DP:128:GLU:O	38:DP:132:LYS:HG3	2.19	0.41
28:DF:114:ILE:HB	28:DF:117:PHE:HB2	2.02	0.41
1:CA:683:G:C5	1:CA:684:A:N7	2.88	0.41
4:CD:75:PHE:HE2	4:CD:207:TYR:HE1	1.67	0.41
23:BA:1374:G:C5	23:BA:1375:C:C5	3.08	0.41
1:CA:625:G:H2'	1:CA:626:U:C6	2.55	0.41
23:BA:320:A:H5''	23:BA:321:G:OP1	2.19	0.41
36:BN:94:TYR:O	36:BN:117:VAL:HG12	2.20	0.41
39:DQ:28:ARG:CG	39:DQ:38:THR:OG1	2.68	0.41
41:DS:12:ILE:HD12	41:DS:46:PHE:HE2	1.85	0.41
23:DA:2230:G:H1'	46:DX:45:ASN:HB2	2.02	0.41
15:CO:50:HIS:O	15:CO:53:HIS:HB3	2.19	0.41
23:DA:2846:G:C5	23:DA:2847:U:C5	3.08	0.41
18:AR:45:SER:HB3	18:AR:51:LEU:HD21	2.01	0.41
36:BN:100:LEU:HD23	36:BN:112:ALA:HA	2.02	0.41
23:BA:1851:U:H2'	23:BA:1852:C:H6	1.85	0.41
23:BA:1284:A:H2'	23:BA:1285:G:O4'	2.20	0.41
5:AE:41:VAL:CG1	5:AE:113:ALA:HA	2.49	0.41
23:DA:1284:A:H2'	23:DA:1285:G:O4'	2.19	0.41
1:AA:35:G:C6	1:AA:36:C:N4	2.87	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:CB:140:HIS:O	2:CB:144:ARG:HG2	2.20	0.41
33:DK:90:GLN:O	33:DK:91:LEU:HB2	2.20	0.41
1:CA:1448:C:H2'	1:CA:1449:C:H6	1.84	0.41
35:DM:121:ALA:HA	35:DM:124:LYS:HG3	2.02	0.41
23:BA:2718:G:H2'	23:BA:2719:G:C8	2.55	0.41
23:DA:725:G:C6	23:DA:726:G:N1	2.88	0.41
23:BA:2338:G:C2	23:BA:2339:G:C8	3.08	0.41
1:CA:697:U:H2'	1:CA:698:G:H5'	2.01	0.41
27:BE:117:ARG:HD2	27:BE:190:GLU:O	2.20	0.41
1:AA:119:A:C5	1:AA:240:C:C4	3.08	0.41
43:DU:95:LYS:HB3	43:DU:99:CYS:O	2.19	0.41
23:BA:71:A:H4'	23:BA:72:U:H5''	2.02	0.41
37:BO:66:ALA:HA	37:BO:69:VAL:HG12	2.02	0.41
46:DX:11:ARG:HH11	46:DX:60:PHE:HA	1.85	0.41
46:BX:11:ARG:HG3	46:BX:62:VAL:HA	2.02	0.41
25:DC:27:THR:CG2	25:DC:83:GLU:HG2	2.50	0.41
34:BL:57:THR:HG23	34:BL:59:LEU:HB3	1.99	0.41
23:BA:2210:G:C3'	23:BA:2210:G:N3	2.77	0.41
36:DN:14:SER:HA	36:DN:17:ARG:HG2	2.02	0.41
23:DA:342:G:O2'	23:DA:343:C:H5''	2.20	0.41
43:BU:17:SER:CB	43:BU:71:LYS:HD2	2.50	0.41
23:BA:2287:A:C5	23:BA:2289:G:C5	3.08	0.41
23:DA:1518:C:O2'	23:DA:1519:G:H5'	2.20	0.41
4:AD:30:LYS:C	4:AD:32:ALA:N	2.70	0.41
44:DV:179:ASP:CG	44:DV:180:VAL:N	2.74	0.41
23:BA:388:G:C4	23:BA:390:A:C6	3.09	0.41
1:AA:1368:G:O2'	1:AA:1369:C:H5'	2.20	0.41
44:DV:24:LEU:HD21	44:DV:86:VAL:HG23	2.01	0.41
23:BA:1486:A:N6	23:BA:1504:C:N4	2.68	0.41
5:CE:48:ALA:HA	5:CE:49:PRO:HD3	1.91	0.41
23:DA:1388:G:N3	23:DA:1389:G:C8	2.89	0.41
35:BM:127:ILE:HG22	35:BM:128:LYS:N	2.36	0.41
23:DA:2849:U:H4'	23:DA:2868:A:C2	2.55	0.41
23:DA:2809:A:N1	23:DA:2892:A:C4	2.88	0.41
1:AA:841:U:O2'	1:AA:842:C:H6	2.03	0.41
23:DA:828:U:C5	23:DA:829:A:N6	2.88	0.41
23:BA:2306:C:H4'	28:BF:136:ARG:NH2	2.34	0.41
25:BC:126:GLN:HG2	25:BC:127:VAL:H	1.85	0.41
44:DV:9:TYR:O	44:DV:38:TYR:HE2	2.03	0.41
22:AV:6165:G:C6	22:AV:6166:U:C4	3.08	0.41
1:AA:1116:C:H3'	1:AA:1117:G:H5''	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:AD:79:PHE:CD1	4:AD:207:TYR:HD1	2.38	0.41
1:CA:626:U:H5''	16:CP:38:TYR:CD2	2.55	0.41
13:AM:19:LEU:HA	13:AM:22:ILE:HG12	2.03	0.41
1:AA:176:C:H5''	20:AT:29:LYS:HZ1	1.85	0.41
20:AT:33:ILE:CD1	20:AT:62:LEU:HD22	2.50	0.41
1:AA:17:U:N3	1:AA:18:C:C4	2.88	0.41
23:BA:1027:A:N6	23:BA:1126:A:N9	2.67	0.41
1:CA:1289:A:C8	1:CA:1290:G:C8	3.08	0.41
21:AU:6:ARG:HG3	21:AU:15:ARG:HH11	1.85	0.41
23:DA:1926:U:O2	23:DA:1929:G:C2	2.73	0.41
1:AA:1289:A:C8	1:AA:1290:G:C8	3.08	0.41
25:BC:105:ILE:HG12	25:BC:106:ILE:HD12	2.01	0.41
38:BP:110:ILE:HD12	38:BP:110:ILE:HA	1.86	0.41
32:BJ:27:TYR:O	32:BJ:29:PRO:HD3	2.20	0.41
23:BA:2723:C:H6	23:BA:2723:C:O5'	2.03	0.41
23:DA:1465:G:C2	23:DA:1466:G:C8	3.08	0.41
1:CA:173:U:C6	1:CA:197:A:C2	3.08	0.41
25:BC:202:LYS:HG3	25:BC:203:ASN:OD1	2.19	0.41
23:BA:735:A:H3'	23:BA:736:C:C6	2.54	0.41
3:CC:134:ILE:HD11	3:CC:153:VAL:HG22	2.01	0.41
23:DA:2194:G:H2'	23:DA:2195:C:H6	1.85	0.41
37:DO:102:ALA:HA	37:DO:105:ALA:HB3	2.02	0.41
1:CA:35:G:C6	1:CA:36:C:N4	2.88	0.41
1:AA:584:G:H2'	1:AA:585:G:C8	2.55	0.41
29:BG:144:VAL:O	29:BG:148:ILE:HG12	2.20	0.41
3:AC:108:ASN:HA	3:AC:109:PRO:HD2	1.87	0.41
23:BA:2509:G:C5	23:BA:2510:C:C5	3.08	0.41
23:BA:2065:C:H2'	23:BA:2066:C:C6	2.55	0.41
23:DA:2518:A:H5'	23:DA:2518:A:C8	2.55	0.41
23:DA:1502:C:H6	23:DA:1502:C:O5'	2.04	0.41
44:BV:157:LEU:HD12	44:BV:157:LEU:N	2.35	0.41
1:CA:786:G:H2'	1:CA:787:A:O4'	2.21	0.41
23:DA:2304:G:H1	23:DA:2312:U:H3	1.67	0.41
23:DA:1688:U:O2	23:DA:1700:A:H8	2.03	0.41
23:BA:826:U:H2'	23:BA:828:U:O4'	2.20	0.41
23:DA:2057:A:H2'	23:DA:2058:A:O4'	2.20	0.41
42:DT:24:GLY:HA3	42:DT:82:GLN:HE22	1.85	0.41
43:BU:14:LEU:HD23	43:BU:14:LEU:C	2.40	0.41
28:BF:5:LEU:HD21	49:B1:50:THR:CG2	2.50	0.41
43:BU:2:ARG:HG2	43:BU:3:VAL:N	2.35	0.41
10:AJ:78:ASN:O	10:AJ:82:ILE:HG12	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:940:C:H2'	1:CA:941:G:H8	1.85	0.41
25:BC:32:SER:HA	25:BC:36:PRO:HG2	2.01	0.41
23:DA:125:G:H5'	52:D4:19:ARG:CG	2.50	0.41
7:AG:70:LYS:HG3	7:AG:96:GLN:HB3	2.02	0.41
28:DF:53:LEU:HD13	28:DF:88:ILE:HG12	2.02	0.41
42:BT:15:GLU:N	42:BT:15:GLU:CD	2.62	0.41
23:DA:558:G:P	32:DJ:134:PRO:HD2	2.61	0.41
13:AM:3:ARG:HH12	28:BF:113:ARG:HD2	1.83	0.41
32:DJ:90:LEU:CD1	32:DJ:90:LEU:H	2.30	0.41
35:BM:24:GLY:HA2	35:BM:101:ARG:CA	2.46	0.41
44:DV:102:LEU:HD21	44:DV:124:ILE:HD11	2.01	0.41
23:DA:588:U:H2'	23:DA:589:C:H6	1.81	0.41
23:DA:2686:G:C5	23:DA:2687:U:C4	3.08	0.41
23:DA:1939:U:OP1	23:DA:2604:U:O2'	2.36	0.41
35:BM:20:ALA:O	35:BM:21:THR:O	2.38	0.41
27:BE:111:ALA:HB2	27:BE:206:ILE:HD12	2.00	0.41
23:BA:448:U:H1'	27:BE:84:VAL:HG21	2.02	0.41
23:BA:2593:U:H2'	23:BA:2594:C:H6	1.81	0.41
52:D4:21:ARG:CB	52:D4:31:LEU:HD21	2.50	0.41
23:BA:2476:A:N1	23:BA:2477:C:C4	2.88	0.41
46:DX:23:LYS:HG3	46:DX:23:LYS:O	2.20	0.41
40:BR:28:GLU:HB3	40:BR:29:PRO:HD2	2.02	0.41
38:DP:64:ARG:HA	38:DP:72:VAL:O	2.20	0.41
2:CB:97:TRP:CH2	2:CB:176:GLU:CD	2.93	0.41
16:AP:49:LEU:HD12	16:AP:50:LYS:N	2.35	0.41
11:AK:54:ARG:HG2	11:AK:54:ARG:H	1.67	0.41
4:AD:75:PHE:HE2	4:AD:207:TYR:HE1	1.67	0.41
1:CA:950:U:H4'	1:CA:971:G:N2	2.36	0.41
23:BA:1496:A:C8	23:BA:1577:C:O2'	2.71	0.41
16:CP:47:ASP:C	16:CP:49:LEU:H	2.23	0.41
1:AA:1281:U:O2'	1:AA:1282:C:P	2.78	0.41
23:DA:486:C:H4'	41:DS:60:ASN:ND2	2.35	0.41
10:CJ:30:SER:HB2	10:CJ:80:LYS:HG3	2.02	0.41
23:DA:1028:A:H61	23:DA:1125:G:H2'	1.84	0.41
45:DW:36:ILE:HG23	45:DW:58:THR:HG23	2.03	0.41
23:BA:2013:A:H4'	41:BS:96:ILE:HD12	2.02	0.41
23:BA:1354:A:O5'	23:BA:1354:A:H8	2.03	0.41
23:DA:1464:C:O2	23:DA:1528:A:H2	2.03	0.41
23:DA:1682:G:H2'	23:DA:1683:C:C6	2.55	0.41
1:AA:1020:U:H2'	1:AA:1021:G:C8	2.55	0.41
23:DA:2718:G:H2'	23:DA:2719:G:C8	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:CG:23:VAL:HG13	7:CG:43:PHE:CE2	2.55	0.41
11:AK:34:ASP:HB2	11:AK:35:PRO:HD2	2.02	0.41
23:DA:1328:G:H2'	23:DA:1330:C:C5	2.55	0.41
1:CA:1089:G:C6	1:CA:1090:U:C4	3.08	0.41
2:AB:140:HIS:O	2:AB:144:ARG:HG2	2.19	0.41
23:DA:1998:G:H2'	23:DA:1999:C:C6	2.56	0.41
1:AA:123:C:OP1	1:AA:312:C:H5'	2.20	0.41
23:BA:1676:A:N6	23:BA:1677:A:C6	2.88	0.41
23:DA:396:G:H1'	46:DX:42:GLN:OE1	2.20	0.41
23:BA:2070:G:H2'	23:BA:2071:A:O4'	2.20	0.41
34:BL:67:MET:HA	34:BL:67:MET:HE3	2.02	0.41
38:BP:115:ARG:HG2	38:BP:115:ARG:H	1.51	0.41
5:AE:15:ARG:O	5:AE:15:ARG:HG2	2.21	0.41
13:AM:56:LEU:HD13	13:AM:56:LEU:O	2.20	0.41
5:CE:15:ARG:HG2	5:CE:15:ARG:O	2.20	0.41
52:B4:47:ARG:O	52:B4:48:LYS:HB2	2.20	0.41
1:AA:1352:C:H2'	1:AA:1353:G:C8	2.56	0.41
2:AB:167:PRO:O	2:AB:171:ALA:HB2	2.20	0.41
34:DL:36:LYS:O	34:DL:38:GLN:HG2	2.20	0.41
39:BQ:106:PHE:HA	39:BQ:109:LEU:HD12	2.02	0.41
32:BJ:88:LYS:O	32:BJ:92:GLN:HB3	2.20	0.41
34:BL:62:LEU:HA	34:BL:63:PRO:HD3	1.77	0.41
23:BA:960:A:H61	35:BM:82:ARG:NH2	2.18	0.41
44:DV:97:GLU:O	44:DV:98:MET:HB3	2.21	0.41
10:AJ:33:GLN:HB2	10:AJ:75:ILE:CD1	2.50	0.41
23:DA:572:A:H2'	23:DA:573:G:O4'	2.20	0.41
3:AC:59:ARG:HA	3:AC:63:ASN:O	2.20	0.41
23:DA:2542:A:H8	23:DA:2544:G:O6	2.03	0.41
3:CC:22:TRP:HE3	3:CC:23:TYR:O	2.02	0.41
27:BE:181:LEU:CD2	27:BE:186:ILE:HD11	2.50	0.41
37:DO:31:SER:HB3	37:DO:34:HIS:H	1.84	0.41
13:AM:3:ARG:HH11	28:BF:113:ARG:HH11	1.69	0.41
23:BA:796:C:H2'	23:BA:797:C:H6	1.81	0.41
11:AK:91:ARG:O	11:AK:95:ILE:HG13	2.20	0.41
28:DF:122:PRO:O	28:DF:125:PHE:HD1	2.03	0.41
11:CK:91:ARG:O	11:CK:95:ILE:HG13	2.19	0.41
46:DX:84:GLY:O	46:DX:85:LEU:C	2.58	0.41
32:BJ:114:LEU:HA	32:BJ:118:PRO:HB3	2.01	0.41
25:BC:69:ARG:HH12	25:BC:117:VAL:HG21	1.85	0.41
27:BE:39:TRP:CD1	27:BE:101:LEU:HB2	2.56	0.41
23:BA:595:C:H2'	23:BA:596:G:O4'	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1422:G:H5'	33:DK:48:PRO:HB3	2.01	0.41
38:BP:100:TYR:C	38:BP:102:ILE:N	2.74	0.41
1:CA:57:G:C6	1:CA:58:C:C4	3.09	0.41
28:DF:115:ARG:HD2	28:DF:115:ARG:N	2.35	0.41
1:CA:1038:C:H2'	1:CA:1039:C:H6	1.85	0.41
38:BP:77:PRO:HB2	38:BP:80:SER:HB2	2.03	0.41
13:CM:8:GLU:OE1	13:CM:22:ILE:HG23	2.21	0.41
39:BQ:28:ARG:HG3	39:BQ:38:THR:OG1	2.19	0.41
4:CD:138:TYR:CD1	4:CD:138:TYR:C	2.94	0.41
23:BA:486:C:H4'	41:BS:60:ASN:ND2	2.35	0.41
23:DA:2050:C:H1'	26:DD:156:MET:CE	2.49	0.41
23:BA:1289:C:H2'	23:BA:1290:C:C6	2.55	0.41
45:BW:14:ARG:HB2	45:BW:14:ARG:HE	1.50	0.41
1:AA:1057:G:H2'	1:AA:1058:G:O4'	2.20	0.41
23:BA:2783:G:N2	26:BD:37:ARG:HH12	2.19	0.41
23:DA:1414:G:C5	23:DA:1415:U:C5	3.08	0.41
4:CD:78:LEU:O	4:CD:81:GLU:HB3	2.21	0.41
1:AA:659:U:O2'	1:AA:660:G:H5'	2.20	0.41
2:CB:167:PRO:O	2:CB:171:ALA:HB2	2.20	0.41
23:BA:2228:G:C6	23:BA:2229:C:C4	3.09	0.41
23:BA:1203:G:C6	23:BA:1204:A:C6	3.09	0.41
20:AT:82:SER:O	20:AT:86:ARG:HB3	2.21	0.41
23:DA:2029:G:H2'	23:DA:2031:A:OP1	2.20	0.41
35:DM:18:LYS:HB3	35:DM:19:GLY:H	1.62	0.41
23:DA:1396:U:O2	23:DA:1396:U:H2'	2.18	0.41
1:AA:1432:G:O5'	1:AA:1432:G:H8	2.02	0.41
34:BL:81:GLN:HE21	34:BL:81:GLN:HB2	1.59	0.41
48:BZ:55:ARG:HD3	48:BZ:55:ARG:HA	1.65	0.41
17:AQ:11:VAL:O	17:AQ:11:VAL:HG22	2.21	0.41
23:BA:312:G:H5'	23:BA:331:A:H2'	2.01	0.41
23:DA:1214:A:H2'	23:DA:1215:G:O4'	2.20	0.41
23:BA:1695:G:N2	23:BA:1696:G:C8	2.88	0.41
37:BO:102:ALA:HA	37:BO:105:ALA:HB3	2.03	0.41
1:AA:366:C:O2'	1:AA:367:U:P	2.78	0.41
23:BA:1140:C:OP1	32:BJ:46:LEU:HB3	2.21	0.41
23:BA:2420:C:P	53:B5:34:TRP:HA	2.60	0.41
23:DA:1022:G:C5	23:DA:1140:C:N4	2.88	0.41
28:DF:5:LEU:HD23	28:DF:5:LEU:HA	1.84	0.41
29:BG:86:GLU:HG2	29:BG:164:TYR:O	2.19	0.41
42:DT:30:VAL:HG21	42:DT:79:ALA:CB	2.50	0.41
50:D2:44:THR:HG22	50:D2:45:VAL:N	2.36	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
37:DO:65:VAL:O	37:DO:69:VAL:HG12	2.20	0.41
1:AA:940:C:H2'	1:AA:941:G:H8	1.86	0.41
10:CJ:78:ASN:O	10:CJ:82:ILE:HG12	2.20	0.41
25:DC:136:ILE:HA	25:DC:137:PRO:HD3	1.96	0.41
1:CA:1372:U:OP1	9:CI:71:SER:HB3	2.21	0.41
23:DA:971:C:H2'	23:DA:972:G:H5'	2.03	0.41
25:BC:35:LYS:CB	25:BC:36:PRO:HD3	2.50	0.41
25:DC:61:LEU:HD12	25:DC:61:LEU:HA	1.84	0.41
23:BA:1493:C:C4	23:BA:2210:G:O2'	2.73	0.41
23:DA:330:A:O2'	23:DA:331:A:H8	2.04	0.41
23:BA:342:G:O2'	23:BA:343:C:H5''	2.21	0.41
32:BJ:57:LEU:HA	32:BJ:57:LEU:HD13	1.72	0.41
43:DU:17:SER:CB	43:DU:71:LYS:HD2	2.50	0.41
26:BD:181:LEU:HD13	26:BD:181:LEU:HA	1.73	0.41
23:DA:1543:A:H3'	23:DA:1543:A:H8	1.85	0.41
28:DF:171:ALA:O	28:DF:175:LEU:HG	2.20	0.41
27:DE:155:LEU:HA	27:DE:174:VAL:HG23	2.03	0.41
23:BA:1543:A:H8	23:BA:1543:A:H3'	1.85	0.41
23:DA:1612:C:O3'	52:D4:5:TRP:HD1	2.03	0.41
1:CA:1428:A:H2'	1:CA:1429:C:C6	2.55	0.41
23:BA:482:A:H8	23:BA:482:A:O5'	2.04	0.41
37:BO:61:ASN:O	37:BO:62:LYS:C	2.59	0.41
29:DG:105:LEU:HD22	29:DG:113:VAL:HB	2.02	0.41
38:DP:100:TYR:C	38:DP:102:ILE:N	2.73	0.41
4:CD:63:LYS:O	4:CD:67:ILE:HG13	2.21	0.41
23:DA:1951:U:O2	23:DA:1953:A:H8	2.02	0.41
23:DA:1276:A:H5''	23:DA:1276:A:H8	1.85	0.41
25:DC:108:PRO:HG3	25:DC:143:HIS:NE2	2.35	0.41
34:BL:126:VAL:HG22	34:BL:145:PRO:HB2	2.02	0.41
23:BA:1265:A:H3'	50:B2:19:ARG:NH1	2.36	0.41
35:DM:104:PHE:HE1	35:DM:125:LEU:HD11	1.84	0.41
1:AA:1038:C:H2'	1:AA:1039:C:H6	1.84	0.41
23:DA:302:C:H2'	23:DA:303:U:C6	2.56	0.41
23:DA:302:C:O2'	23:DA:303:U:H5'	2.21	0.41
23:DA:1494:A:H4'	23:DA:1495:A:OP1	2.21	0.41
4:CD:90:GLY:HA2	4:CD:204:ILE:HD11	2.02	0.41
32:DJ:151:HIS:CD2	32:DJ:152:PRO:O	2.73	0.41
35:BM:61:GLY:O	44:BV:177:PRO:HA	2.20	0.41
44:DV:38:TYR:CD1	44:DV:38:TYR:O	2.73	0.41
22:CV:6165:G:C6	22:CV:6166:U:C4	3.09	0.41
9:AI:86:VAL:CG2	9:AI:93:ARG:HB2	2.51	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BA:2487:G:H2'	23:BA:2488:A:C8	2.56	0.41
4:AD:138:TYR:C	4:AD:138:TYR:CD1	2.93	0.41
23:DA:719:C:H2'	23:DA:720:C:C6	2.53	0.41
23:BA:755:C:H2'	23:BA:756:C:C6	2.55	0.41
23:BA:1270:C:H5''	23:BA:1271:G:C5'	2.50	0.41
23:BA:699:A:H2'	23:BA:700:G:O4'	2.20	0.41
4:CD:188:LEU:HD12	4:CD:188:LEU:H	1.85	0.41
1:CA:698:G:C6	1:CA:699:C:C4	3.08	0.41
23:BA:1696:G:C6	23:BA:1697:G:C4	3.08	0.41
23:DA:1109:C:H2'	23:DA:1110:G:O4'	2.20	0.41
44:BV:161:VAL:HG12	44:BV:162:GLU:N	2.36	0.41
4:AD:33:MET:HG2	4:AD:37:PRO:HA	2.03	0.41
23:DA:1170:G:N2	23:DA:1180:C:C2	2.89	0.41
23:DA:30:G:C5	23:DA:31:C:C4	3.09	0.41
23:BA:1581:G:H2'	23:BA:1582:C:O4'	2.20	0.41
17:CQ:11:VAL:O	17:CQ:11:VAL:HG22	2.21	0.41
23:BA:2284:C:O5'	23:BA:2284:C:H6	2.03	0.41
1:AA:743:U:O5'	1:AA:743:U:H6	2.04	0.41
48:BZ:52:HIS:CD2	48:BZ:52:HIS:H	2.37	0.41
6:CF:70:ASP:OD1	6:CF:70:ASP:N	2.51	0.41
32:BJ:101:TYR:N	32:BJ:101:TYR:CD1	2.88	0.41
1:CA:290:C:O5'	1:CA:290:C:H6	2.03	0.41
14:CN:40:CYS:SG	14:CN:42:ILE:HB	2.60	0.41
7:CG:25:ALA:O	7:CG:29:LYS:HG2	2.20	0.41
34:BL:36:LYS:O	34:BL:38:GLN:HG2	2.19	0.41
23:BA:2057:A:H2'	23:BA:2058:A:O4'	2.21	0.41
40:DR:22:VAL:CG1	40:DR:23:GLU:H	2.33	0.41
40:BR:38:LEU:O	40:BR:52:VAL:HG12	2.21	0.41
23:BA:114(B):A:C2	23:BA:1144:G:C8	3.08	0.41
23:DA:2416:C:OP1	34:DL:64:LYS:O	2.39	0.41
23:DA:2422:A:C5	23:DA:2424:C:C4	3.09	0.41
34:BL:85:LEU:HA	34:BL:88:LEU:HB2	2.01	0.41
23:BA:675:A:C4'	27:BE:67:GLN:NE2	2.81	0.41
43:BU:95:LYS:HB3	43:BU:99:CYS:O	2.20	0.41
47:BY:6:VAL:O	47:BY:9:GLN:HB2	2.21	0.41
46:DX:11:ARG:HG3	46:DX:62:VAL:HA	2.02	0.41
23:BA:846:C:H4'	23:BA:847:U:O5'	2.21	0.41
23:BA:933:A:H2'	23:BA:934:G:H5'	2.03	0.41
1:AA:674:G:P	6:AF:87:ARG:HH22	2.42	0.41
2:AB:80:ILE:HD11	2:AB:208:ILE:CG2	2.51	0.41
30:BH:133:HIS:HD2	30:BH:135:GLU:HG2	1.85	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
28:BF:87:PRO:O	28:BF:88:ILE:HB	2.21	0.41
33:BK:71:ARG:NH2	33:BK:77:ILE:HG21	2.35	0.41
30:DH:82:ARG:C	30:DH:89:TYR:HB2	2.41	0.41
24:DB:72:G:N2	24:DB:103:U:C5	2.88	0.41
23:BA:643:A:OP1	51:B3:42:TRP:NE1	2.54	0.41
28:DF:15:VAL:HG22	28:DF:175:LEU:HB3	2.03	0.41
28:DF:32:PRO:CB	28:DF:172:LEU:HD22	2.49	0.41
28:BF:122:PRO:O	28:BF:125:PHE:HD1	2.02	0.41
51:D3:34:LEU:HD13	51:D3:34:LEU:H	1.86	0.41
23:DA:477:A:H2'	23:DA:478:A:C8	2.54	0.41
22:AV:6171:U:O2'	22:AV:6172:U:P	2.79	0.41
2:CB:71:VAL:HG23	2:CB:164:VAL:HG13	2.02	0.41
23:DA:2777:G:C4'	23:DA:2778:A:H5'	2.50	0.41
34:BL:9:ASN:N	34:BL:10:PRO:CD	2.83	0.41
1:AA:370:C:O2'	1:AA:371:G:H5'	2.21	0.41
35:DM:127:ILE:HG22	35:DM:128:LYS:N	2.35	0.41
12:CL:51:LEU:HD12	12:CL:51:LEU:N	2.34	0.41
27:DE:150:GLY:HA2	27:DE:172:TRP:CD2	2.55	0.41
27:BE:150:GLY:HA2	27:BE:172:TRP:CD2	2.56	0.41
23:DA:1952:A:C2	33:DK:22:ILE:HG23	2.56	0.41
28:DF:128:ARG:HH21	28:DF:130:ASN:ND2	2.17	0.41
23:BA:302:C:H2'	23:BA:303:U:C6	2.55	0.41
1:AA:817:C:H1'	1:AA:819:A:C5'	2.51	0.41
35:DM:26:TYR:CD1	35:DM:26:TYR:O	2.73	0.41
24:BB:48:A:H2'	24:BB:49:C:C6	2.55	0.41
20:AT:63:ILE:HG21	20:AT:81:LYS:HG3	2.02	0.41
1:CA:865:A:H5'	1:CA:1078:U:C4	2.56	0.41
23:DA:2257:U:O2'	23:DA:2258:C:H5'	2.21	0.41
1:AA:1229:A:H2'	1:AA:1230:C:H6	1.86	0.41
25:DC:37:LEU:HD12	25:DC:38:LYS:N	2.36	0.41
23:BA:2341:G:H2'	23:BA:2342:C:C6	2.55	0.41
23:BA:1750:G:H2'	23:BA:1751:C:H6	1.84	0.41
1:CA:562:C:H6	1:CA:562:C:H5'	1.85	0.41
27:BE:72:ARG:O	27:BE:73:ALA:O	2.38	0.41
29:DG:83:TYR:CZ	29:DG:138:LYS:HG3	2.55	0.41
23:BA:1351:C:C2	23:BA:1381:G:C2	3.08	0.41
4:AD:156:GLU:O	4:AD:160:GLN:HG3	2.21	0.41
23:DA:1232:G:H2'	23:DA:1233:C:H6	1.85	0.41
27:DE:6:MET:HB3	27:DE:7:TYR:H	1.55	0.41
23:DA:979:G:H3'	23:DA:980:A:H5''	2.02	0.41
7:CG:70:LYS:CG	7:CG:96:GLN:HB3	2.51	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1124:G:H5'	10:AJ:35:SER:HB2	2.01	0.41
23:DA:1818:U:H2'	25:DC:157:ARG:HG3	2.02	0.41
6:CF:91:VAL:HG12	6:CF:92:LYS:O	2.20	0.41
24:BB:52:A:H2'	24:BB:53:A:O4'	2.21	0.41
24:BB:3:C:H2'	24:BB:4:C:C6	2.55	0.41
23:BA:2235:G:H2'	23:BA:2236:C:C6	2.55	0.41
1:AA:407:G:H4'	4:AD:116:GLN:HA	2.02	0.41
1:CA:755:G:H2'	1:CA:756:C:C6	2.56	0.41
28:BF:178:PHE:HA	28:BF:179:PRO:HD3	1.79	0.41
1:CA:1403:C:O5'	1:CA:1403:C:H6	2.03	0.41
36:DN:99:LYS:HB3	36:DN:99:LYS:HE3	1.86	0.41
23:DA:1676:A:N6	23:DA:1677:A:C6	2.88	0.41
1:CA:261:U:H5	20:CT:79:ARG:CZ	2.33	0.41
23:BA:2642:G:O2'	23:BA:2643:G:H5'	2.21	0.41
43:DU:97:ARG:O	43:DU:97:ARG:HG2	2.21	0.41
23:DA:956:G:H22	23:DA:959:A:H3'	1.85	0.41
41:DS:24:ILE:CG2	41:DS:36:LEU:HD21	2.39	0.41
3:AC:70:VAL:O	3:AC:106:VAL:HG23	2.19	0.41
23:DA:846:C:H4'	23:DA:847:U:O5'	2.20	0.41
23:BA:2721:A:H1'	23:BA:2873:A:O2'	2.21	0.41
48:BZ:26:LEU:HD21	48:BZ:46:ASN:HB3	2.01	0.41
34:BL:6:LEU:N	34:BL:6:LEU:HD23	2.25	0.41
23:DA:773:U:C5'	25:DC:47:GLY:HA3	2.51	0.41
26:BD:171:GLU:HG2	26:BD:185:LYS:CG	2.50	0.41
12:CL:69:ILE:HD12	12:CL:69:ILE:N	2.36	0.41
23:BA:1407:C:H2'	23:BA:1408:C:C6	2.55	0.41
38:BP:74:ARG:HD3	38:BP:76:PHE:CE2	2.55	0.41
36:BN:9:LYS:C	36:BN:10:LEU:CG	2.89	0.41
23:DA:310:A:OP1	43:DU:17:SER:O	2.38	0.41
30:BH:109:ILE:HD13	30:BH:109:ILE:N	2.35	0.41
28:BF:161:THR:HG21	28:BF:172:LEU:CD2	2.50	0.41
28:DF:13:GLU:O	28:DF:14:GLU:HB2	2.19	0.41
3:AC:34:LEU:HD21	3:AC:38:ARG:HH21	1.84	0.41
24:DB:106:G:C5	24:DB:107:U:C5	3.08	0.41
4:CD:104:VAL:O	4:CD:108:LEU:HB2	2.21	0.41
34:DL:7:ARG:O	34:DL:10:PRO:HD3	2.21	0.41
9:AI:29:ASN:ND2	9:AI:65:VAL:O	2.51	0.41
8:CH:64:LYS:CG	8:CH:79:VAL:HG21	2.49	0.41
2:CB:8:LYS:HA	2:CB:217:ARG:NH1	2.33	0.41
9:AI:112:LYS:HA	9:AI:119:ALA:HB2	2.02	0.41
23:BA:1308:A:H2'	23:BA:1309:G:O4'	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BA:1804:C:H6	23:BA:1804:C:O5'	2.04	0.41
23:DA:426:C:C2	23:DA:427:U:C6	3.09	0.41
27:DE:12:LEU:HD13	27:DE:17:ARG:HG2	2.03	0.41
23:DA:1496:A:C8	23:DA:1498:C:N3	2.88	0.41
1:CA:818:G:H1'	1:CA:820:U:H5	1.85	0.41
23:BA:1495:A:N3	23:BA:1496:A:C2	2.89	0.41
38:DP:3:ARG:HD2	38:DP:6:LEU:HD23	2.03	0.41
1:AA:1053:G:C3'	1:AA:1054:C:H5'	2.51	0.41
8:AH:11:THR:HA	8:AH:14:ARG:HH12	1.85	0.41
23:DA:2296:U:O4	37:DO:13:ARG:NH2	2.50	0.41
2:AB:22:LYS:HZ2	2:AB:22:LYS:HA	1.86	0.41
41:BS:84:ARG:HB2	41:BS:96:ILE:CG2	2.51	0.41
1:CA:832:C:O2'	1:CA:833:U:H6	2.03	0.41
23:DA:1050:A:H2'	23:DA:1051:G:C8	2.55	0.41
23:BA:1854:A:H62	23:BA:1888:G:H8	1.69	0.41
25:DC:105:ILE:HG13	25:DC:106:ILE:HD12	2.02	0.41
23:DA:1798:U:C5'	25:DC:259:THR:O	2.69	0.41
13:AM:86:CYS:SG	13:AM:88:ARG:HB2	2.61	0.41
1:AA:336:C:H2'	1:AA:337:C:C6	2.55	0.41
1:AA:1058:G:H2'	1:AA:1059:C:C6	2.55	0.41
7:CG:91:VAL:HG12	7:CG:95:ARG:HB3	2.03	0.41
1:CA:1010:G:H2'	1:CA:1011:G:H8	1.86	0.41
23:BA:2329:G:H2'	23:BA:2330:G:C8	2.56	0.41
5:CE:41:VAL:CG1	5:CE:113:ALA:HA	2.51	0.41
23:BA:1414:G:C5	23:BA:1415:U:C5	3.09	0.41
23:DA:335:C:H2'	23:DA:336:C:C6	2.55	0.41
23:DA:1676:A:H2'	23:DA:1677:A:O4'	2.20	0.41
46:BX:70:VAL:O	46:BX:74:VAL:HG23	2.21	0.41
39:BQ:66:ASN:HD21	39:BQ:70:ARG:HH21	1.69	0.41
23:BA:2692:C:H2'	23:BA:2693:A:H8	1.86	0.41
7:CG:80:VAL:C	7:CG:82:GLY:H	2.24	0.41
23:DA:1248:G:C5	39:DQ:3:ARG:HB2	2.55	0.41
1:AA:1408:A:C6	1:AA:1494:G:C6	3.09	0.41
23:DA:1648:C:H2'	23:DA:1649:G:O5'	2.20	0.41
23:BA:1214:A:H2'	23:BA:1215:G:O4'	2.21	0.41
23:BA:2618:G:C6	23:BA:2619:C:C4	3.09	0.41
27:BE:78:ILE:HD12	27:BE:78:ILE:H	1.85	0.41
4:AD:179:GLU:CD	4:AD:179:GLU:H	2.24	0.41
30:BH:26:ALA:HA	30:BH:30:LEU:HB2	2.02	0.41
23:BA:1794:U:H1'	23:BA:1900:A:N3	2.35	0.41
42:BT:24:GLY:HA3	42:BT:82:GLN:HE22	1.86	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
28:DF:66:GLN:HG2	28:DF:67:LYS:N	2.26	0.41
29:BG:101:ARG:HB2	29:BG:117:PRO:HG2	2.03	0.41
43:BU:76:CYS:CB	43:BU:96:ILE:HD13	2.50	0.41
10:CJ:48:THR:HG22	10:CJ:62:HIS:ND1	2.35	0.41
28:DF:41:GLN:HB3	28:DF:43:LEU:HD13	2.03	0.41
28:BF:41:GLN:HB3	28:BF:43:LEU:HD13	2.03	0.41
44:BV:71:VAL:HG11	44:BV:74:VAL:HG23	2.02	0.41
53:D5:13:ARG:HG3	53:D5:14:VAL:HG23	2.02	0.41
1:CA:1371:G:C6	1:CA:1372:U:C4	3.09	0.41
37:DO:17:ARG:HG2	37:DO:18:ILE:HD13	2.02	0.41
37:BO:100:ALA:HA	37:BO:103:GLU:HB3	2.01	0.41
23:DA:1447:G:N3	23:DA:1545:A:H2	2.18	0.41
23:DA:861:A:C2	23:DA:917:A:C4	3.09	0.41
23:BA:1543:A:H3'	23:BA:1543:A:C8	2.56	0.41
23:BA:861:A:C2	23:BA:917:A:C4	3.09	0.41
41:DS:14:PRO:O	41:DS:18:ARG:HG3	2.21	0.41
32:BJ:116:THR:HG23	32:BJ:117:HIS:H	1.86	0.41
1:CA:878:G:C5'	8:CH:89:PRO:HG2	2.46	0.41
10:CJ:54:PHE:HB3	10:CJ:55:LYS:H	1.63	0.41
35:DM:20:ALA:O	35:DM:21:THR:O	2.39	0.41
27:BE:101:LEU:O	27:BE:106:ARG:NH1	2.53	0.41
1:AA:754:C:O2	1:AA:754:C:C3'	2.67	0.41
23:BA:2850:A:H2'	23:BA:2851:A:O4'	2.20	0.41
23:DA:1257:C:H4'	27:DE:83:PHE:CD2	2.56	0.41
23:BA:85:G:C8	23:BA:85:G:H5''	2.51	0.41
10:CJ:6:ILE:HB	10:CJ:98:ILE:HG12	2.02	0.41
23:DA:2306:C:H4'	28:DF:136:ARG:NH2	2.35	0.41
1:CA:429:U:H4'	1:CA:430:A:O5'	2.20	0.41
46:BX:37:ILE:HG23	46:BX:38:SER:N	2.36	0.41
23:BA:1952:A:C2	33:BK:22:ILE:HG23	2.55	0.41
18:CR:35:ARG:O	18:CR:37:VAL:N	2.49	0.41
1:AA:815:A:H4'	1:AA:817:C:C4	2.55	0.41
4:AD:61:LYS:HE3	4:AD:207:TYR:OH	2.20	0.41
1:CA:161:A:H2'	1:CA:162:A:H8	1.86	0.41
1:CA:983:A:H2	1:CA:984:C:C6	2.39	0.41
23:BA:1206:G:C5	23:BA:1207:C:C5	3.09	0.41
1:AA:824:C:C6	1:AA:824:C:C3'	3.04	0.41
52:B4:24:THR:HG23	52:B4:27:GLY:H	1.85	0.41
1:AA:349:A:O2'	1:AA:350:G:H5'	2.21	0.41
23:DA:161:U:O2	23:DA:165:U:O4	2.39	0.41
23:DA:288:C:O2'	23:DA:289:A:H5'	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BA:161:U:O2	23:BA:165:U:O4	2.39	0.41
26:DD:76:ARG:HG2	26:DD:77:ILE:HG13	2.03	0.41
18:CR:45:SER:HB3	18:CR:51:LEU:HG	2.02	0.41
23:BA:2817:G:H21	23:BA:2836:U:C1'	2.33	0.41
23:BA:912:C:C2	23:BA:913:U:C5	3.09	0.41
1:AA:1442:G:H3'	1:AA:1442:G:C8	2.55	0.41
11:CK:41:THR:HG21	11:CK:71:LYS:HB2	2.03	0.41
23:DA:1271:G:O3'	23:DA:1272:A:H4'	2.21	0.41
1:CA:342:C:C2'	1:CA:343:U:H5'	2.51	0.41
29:BG:105:LEU:HD22	29:BG:113:VAL:HB	2.01	0.41
9:AI:46:ALA:O	9:AI:49:PRO:HD2	2.21	0.41
35:DM:34:LEU:HD12	35:DM:130:LYS:O	2.21	0.41
23:BA:1239:G:C6	23:BA:1240:U:C4	3.09	0.41
23:DA:489:G:C5	23:DA:1284:A:C2	3.09	0.41
23:BA:2695:C:H2'	23:BA:2696:U:C6	2.56	0.41
23:DA:1696:G:C6	23:DA:1697:G:C4	3.08	0.41
2:AB:16:HIS:CD2	2:AB:210:SER:HA	2.56	0.41
7:CG:70:LYS:HG3	7:CG:96:GLN:HB3	2.02	0.41
20:AT:94:ALA:C	20:AT:96:GLY:H	2.24	0.41
1:AA:1466:C:H2'	1:AA:1467:G:O4'	2.20	0.41
23:BA:29:U:O2'	23:BA:30:G:H5'	2.21	0.41
42:BT:14:SER:O	42:BT:17:ALA:N	2.54	0.41
1:AA:293:G:C6	1:AA:294:U:C4	3.09	0.41
24:BB:7:G:H5''	37:BO:29:PHE:CD2	2.56	0.41
1:CA:658:G:C6	1:CA:659:U:C4	3.09	0.41
4:CD:199:ASN:ND2	4:CD:202:LEU:HG	2.35	0.41
23:BA:308:G:H2'	23:BA:309:G:O4'	2.21	0.41
24:DB:33:G:O2'	24:DB:34:U:H5'	2.20	0.41
23:DA:1024:G:O5'	23:DA:1024:G:H8	2.03	0.41
24:DB:7:G:H5''	37:DO:29:PHE:CD2	2.55	0.41
44:DV:30:ASN:OD1	44:DV:33:LEU:N	2.53	0.41
1:AA:1089:G:C6	1:AA:1090:U:C4	3.09	0.41
24:DB:116:G:H4'	37:DO:55:ALA:O	2.21	0.41
34:BL:39:LYS:HD3	34:BL:39:LYS:HA	1.71	0.41
34:BL:48:PRO:O	34:BL:49:ARG:C	2.59	0.41
34:DL:47:ASP:HB3	34:DL:48:PRO:HA	2.01	0.41
53:D5:59:LYS:HA	53:D5:62:LEU:HD11	2.03	0.41
23:DA:593:G:C6	23:DA:594:U:C4	3.09	0.41
1:AA:1130:A:H4'	9:AI:20:ARG:HH22	1.86	0.41
40:DR:4:ILE:HG22	40:DR:5:VAL:N	2.36	0.41
23:BA:1825:A:H2'	23:BA:1826:G:C8	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BA:1826:G:C6	23:BA:1827:C:C4	3.09	0.41
23:BA:1164:G:H5''	23:BA:1164:G:H8	1.86	0.41
23:DA:2247:A:O2'	23:DA:2248:C:H5'	2.20	0.41
23:DA:827:U:H2'	23:DA:2068:U:C2	2.56	0.41
23:BA:733:G:N7	23:BA:761:A:N6	2.68	0.41
28:DF:5:LEU:HD21	49:D1:50:THR:CG2	2.50	0.41
1:AA:737:A:H2'	1:AA:738:C:H6	1.85	0.41
3:AC:131:ARG:HH21	5:AE:50:GLU:CG	2.34	0.41
23:DA:71:A:H4'	23:DA:72:U:H5''	2.03	0.41
28:BF:5:LEU:HA	28:BF:5:LEU:HD23	1.85	0.41
37:BO:28:VAL:HG13	37:BO:35:ILE:HD11	2.01	0.41
37:BO:65:VAL:O	37:BO:69:VAL:HG12	2.21	0.41
30:DH:5:LEU:CD2	30:DH:5:LEU:H	2.18	0.41
23:DA:2712:U:O2'	23:DA:2713:A:H5'	2.21	0.41
38:BP:68:TYR:N	38:BP:68:TYR:CD2	2.89	0.41
46:BX:11:ARG:HH11	46:BX:60:PHE:HA	1.85	0.41
23:DA:270(H):C:H2'	23:DA:270(I):C:C6	2.56	0.41
44:BV:51:ALA:HB1	44:BV:57:ILE:HD11	2.02	0.41
44:BV:53:ILE:CG2	44:BV:71:VAL:HB	2.51	0.41
26:BD:116:VAL:HG13	26:BD:117:MET:N	2.35	0.41
36:DN:4:LEU:CG	36:DN:4:LEU:O	2.66	0.41
52:D4:8:ASN:HD21	52:D4:10:ARG:HB3	1.86	0.41
23:DA:2024:G:H2'	23:DA:2025:C:H6	1.86	0.41
23:BA:773:U:C5'	25:BC:47:GLY:HA3	2.51	0.41
25:DC:270:ILE:C	25:DC:271:ILE:HG13	2.40	0.41
23:DA:1006:C:H1'	32:DJ:129:MET:HG2	2.02	0.41
5:AE:20:GLN:O	5:AE:21:ALA:C	2.58	0.41
7:AG:70:LYS:CG	7:AG:96:GLN:HB3	2.51	0.41
23:DA:2636:U:H2'	23:DA:2637:U:H6	1.86	0.41
12:AL:29:ALA:HA	12:AL:30:PRO:HD3	1.80	0.41
1:CA:1329:A:H5''	13:CM:26:GLY:N	2.36	0.41
12:CL:103:VAL:O	12:CL:106:ALA:HB3	2.21	0.41
32:DJ:57:LEU:HA	32:DJ:57:LEU:HD13	1.73	0.41
23:BA:2516:G:C6	23:BA:2517:C:C4	3.08	0.41
43:DU:17:SER:HB2	43:DU:71:LYS:HD2	2.02	0.41
33:BK:47:ILE:HA	33:BK:47:ILE:HD12	1.87	0.41
7:CG:15:ASP:HA	7:CG:24:THR:CG2	2.51	0.41
28:DF:18:GLU:HB3	28:DF:175:LEU:HD13	2.03	0.41
1:AA:392:G:C2	1:AA:393:A:C5	3.09	0.41
27:DE:181:LEU:HA	27:DE:181:LEU:HD23	1.93	0.41
35:DM:24:GLY:HA2	35:DM:100:GLY:C	2.41	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DA:468:G:N7	52:D4:39:ARG:NH2	2.69	0.41
24:BB:106:G:O2'	24:BB:107:U:H5'	2.20	0.41
23:DA:1332:G:N2	23:DA:1609:A:O2'	2.54	0.41
24:DB:74:U:H2'	24:DB:75:G:H8	1.83	0.41
25:DC:235:GLY:C	25:DC:237:GLU:H	2.24	0.41
10:CJ:56:HIS:O	10:CJ:58:ASP:N	2.45	0.41
1:CA:438:G:C4'	1:CA:439:A:OP1	2.67	0.41
27:BE:28:ILE:HA	27:BE:112:MET:HE3	2.03	0.41
1:CA:802:A:H2'	1:CA:803:G:O4'	2.21	0.41
25:BC:108:PRO:HG3	25:BC:143:HIS:NE2	2.36	0.41
23:BA:1258:C:O4'	27:BE:84:VAL:HG11	2.21	0.41
8:AH:64:LYS:CG	8:AH:79:VAL:HG21	2.50	0.41
23:BA:444:C:H2'	23:BA:445:C:H6	1.85	0.41
1:CA:300:A:H3'	1:CA:300:A:C8	2.56	0.41
27:BE:177:ALA:HB1	27:BE:178:PRO:HD2	2.03	0.41
46:DX:37:ILE:HG22	46:DX:38:SER:N	2.36	0.41
23:DA:2476:A:N1	23:DA:2477:C:C4	2.88	0.41
1:AA:502:G:H2'	1:AA:503:C:O4'	2.21	0.41
1:CA:1228:C:H5'	13:CM:115:LYS:O	2.21	0.41
23:BA:332:A:O2'	23:BA:333:G:P	2.79	0.41
9:AI:53:VAL:HG11	9:AI:85:LEU:HD22	2.02	0.41
23:BA:1374:G:H2'	23:BA:1375:C:H6	1.86	0.41
38:DP:118:ARG:HA	38:DP:121:ILE:HD12	2.03	0.41
16:AP:47:ASP:C	16:AP:49:LEU:H	2.25	0.41
23:BA:2749:A:H4'	29:BG:62:LYS:CB	2.51	0.41
1:AA:1116:C:H2'	1:AA:1117:G:H5''	2.03	0.41
7:CG:41:ARG:HB3	7:CG:41:ARG:HH11	1.86	0.41
41:BS:10:VAL:HG21	41:BS:103:ILE:HD13	2.01	0.41
35:BM:132:VAL:HG11	44:BV:81:ARG:NH1	2.36	0.41
45:BW:51:VAL:HG21	45:BW:80:HIS:HA	2.01	0.41
27:DE:40:GLN:O	27:DE:43:LYS:HG2	2.21	0.41
36:DN:104:ARG:HH11	36:DN:104:ARG:CB	2.33	0.41
36:DN:116:LEU:HA	36:DN:116:LEU:HD23	1.79	0.41
36:DN:94:TYR:O	36:DN:117:VAL:HG12	2.20	0.41
35:BM:68:ILE:HD13	35:BM:103:MET:HG3	2.02	0.41
23:DA:1790:C:H5''	23:DA:1791:A:P	2.61	0.41
1:CA:663:A:O2'	1:CA:664:G:H5'	2.21	0.41
40:DR:17:GLY:HA2	40:DR:96:ILE:O	2.21	0.41
23:BA:1798:U:H5''	25:BC:259:THR:O	2.21	0.41
41:DS:22:ASP:HA	41:DS:25:ARG:NH1	2.36	0.41
23:DA:2817:G:H21	23:DA:2836:U:C1'	2.33	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DA:2836:U:H2'	23:DA:2837:G:H8	1.86	0.41
23:DA:1748:G:H2'	23:DA:1749:A:H8	1.84	0.41
23:DA:2765:A:H5'	23:DA:2766:G:OP2	2.21	0.41
23:DA:2658:C:H4'	29:DG:158:HIS:CE1	2.56	0.41
1:CA:1442:G:C8	1:CA:1442:G:H3'	2.55	0.41
1:AA:1272:G:H2'	1:AA:1273:G:H8	1.84	0.41
7:AG:88:PRO:O	7:AG:89:MET:HB3	2.21	0.41
25:DC:140:THR:HG22	25:DC:141:VAL:N	2.36	0.41
23:DA:924:C:H2'	23:DA:925:C:H6	1.86	0.41
23:DA:699:A:H2'	23:DA:700:G:O4'	2.20	0.41
2:AB:166:ASP:HA	2:AB:167:PRO:HD2	1.87	0.41
23:BA:1676:A:O5'	23:BA:1676:A:H8	2.04	0.41
7:CG:92:SER:O	7:CG:96:GLN:HG3	2.21	0.41
11:CK:94:ALA:O	11:CK:98:LEU:HG	2.21	0.41
23:DA:2709:G:O2'	23:DA:2710:C:H5'	2.20	0.41
23:BA:1368:G:C2	23:BA:1369:G:C8	3.09	0.41
29:BG:125:VAL:HG22	29:BG:131:VAL:HG22	2.03	0.41
27:DE:14:PRO:HD3	27:DE:128:ALA:HB2	2.02	0.41
23:BA:2304:G:H1	23:BA:2312:U:H3	1.67	0.41
23:DA:26:G:C6	23:DA:27:G:N1	2.89	0.41
44:BV:155:LEU:HD21	44:BV:171:ILE:HG13	2.02	0.41
27:BE:24:LEU:HA	27:BE:25:PRO:HD3	1.85	0.41
43:DU:34:LYS:HB3	43:DU:34:LYS:HE2	1.80	0.41
23:BA:885:C:O5'	23:BA:885:C:H6	2.04	0.41
50:B2:13:LYS:HE2	50:B2:13:LYS:HB3	1.88	0.41
46:BX:95:LEU:HD22	46:BX:95:LEU:HA	1.82	0.41
24:BB:45:A:H2'	24:BB:45:A:N3	2.35	0.41
23:BA:1765:C:H6	23:BA:1765:C:O5'	2.04	0.41
1:AA:1147:C:O5'	1:AA:1147:C:H6	2.04	0.41
44:DV:43:GLU:O	44:DV:47:VAL:HG23	2.21	0.41
23:BA:533:G:N3	39:BQ:45:TYR:CE1	2.89	0.41
5:CE:106:PRO:O	5:CE:110:LEU:HG	2.21	0.41
23:BA:725:G:C6	23:BA:726:G:N1	2.89	0.41
39:BQ:57:PHE:HB3	39:BQ:61:TRP:CZ2	2.55	0.41
23:DA:997:G:O2'	23:DA:998:C:H5'	2.21	0.41
44:DV:155:LEU:HD21	44:DV:171:ILE:HG13	2.02	0.41
23:BA:1563:G:H2'	23:BA:1564:C:H6	1.85	0.41
23:DA:2019:A:H5''	39:DQ:27:LEU:HD12	2.03	0.41
23:BA:2507:C:C2	23:BA:2508:G:C8	3.09	0.41
1:AA:1051:C:C4	1:AA:1052:U:C4	3.09	0.41
23:DA:447:A:C5	23:DA:473:G:C5	3.09	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
48:DZ:37:LEU:HA	48:DZ:37:LEU:HD23	1.95	0.41
26:DD:141:ILE:O	26:DD:141:ILE:HG13	2.21	0.41
2:CB:36:ARG:N	2:CB:36:ARG:HD2	2.35	0.41
40:BR:75:PHE:CD1	40:BR:75:PHE:C	2.94	0.41
27:BE:6:MET:HB3	27:BE:7:TYR:H	1.55	0.41
12:AL:85:ARG:HH21	12:AL:98:HIS:CG	2.39	0.41
23:DA:583:G:OP2	39:DQ:10:ARG:HD2	2.21	0.41
23:DA:1221:C:H2'	23:DA:122(A):C:C6	2.56	0.41
53:B5:62:LEU:HB2	53:B5:63:PRO:HD3	2.02	0.41
53:B5:59:LYS:HA	53:B5:62:LEU:HD11	2.03	0.41
23:BA:195:A:H4'	23:BA:251:A:O2'	2.21	0.41
40:DR:40:LEU:C	40:DR:45:THR:HB	2.41	0.41
40:BR:22:VAL:CG1	40:BR:23:GLU:H	2.32	0.41
2:CB:69:LEU:HD22	2:CB:159:PRO:HG2	2.02	0.41
28:DF:60:LEU:HD12	28:DF:68:PRO:HB3	2.03	0.41
10:AJ:49:VAL:HG21	14:AN:41:ARG:HB2	2.02	0.41
23:BA:270(K):G:H2'	23:BA:270(L):C:O4'	2.21	0.41
37:DO:66:ALA:HA	37:DO:69:VAL:HG12	2.03	0.41
17:AQ:83:ASP:O	17:AQ:86:GLU:HB2	2.21	0.41
26:BD:117:MET:HB2	26:BD:117:MET:HE3	1.92	0.41
23:DA:948:G:C5'	23:DA:948:G:C8	3.00	0.41
34:BL:57:THR:C	34:BL:59:LEU:N	2.74	0.41
16:CP:4:ILE:H	16:CP:4:ILE:HD12	1.86	0.41
36:DN:14:SER:O	36:DN:15:SER:C	2.59	0.41
28:DF:87:PRO:O	28:DF:88:ILE:HB	2.21	0.41
40:BR:77:ALA:C	40:BR:79:VAL:N	2.73	0.41
19:AS:6:LYS:HD2	19:AS:6:LYS:N	2.34	0.41
48:BZ:40:THR:OG1	48:BZ:43:ILE:HG12	2.20	0.41
26:DD:181:LEU:HD13	26:DD:181:LEU:HA	1.70	0.41
23:DA:1334:G:C6	23:DA:1335:U:C4	3.09	0.41
1:CA:1151:A:O2'	1:CA:1152:A:O5'	2.37	0.41
23:DA:643:A:O2'	23:DA:644:A:H5'	2.21	0.41
23:DA:1778:U:H2'	23:DA:1784:A:H62	1.83	0.41
7:AG:15:ASP:HA	7:AG:24:THR:CG2	2.51	0.41
15:CO:45:VAL:HG23	15:CO:46:HIS:N	2.36	0.41
1:CA:973:G:H4'	10:CJ:54:PHE:O	2.20	0.41
36:DN:100:LEU:HD23	36:DN:112:ALA:HA	2.02	0.41
6:CF:11:ASN:HA	6:CF:12:PRO:HD2	1.86	0.41
13:AM:106:ASN:O	13:AM:107:ALA:CB	2.69	0.41
37:DO:61:ASN:O	37:DO:62:LYS:C	2.59	0.41
27:DE:111:ALA:HB2	27:DE:206:ILE:HD12	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
28:BF:115:ARG:HD2	28:BF:115:ARG:N	2.36	0.41
32:BJ:83:ILE:HG22	32:BJ:84:ARG:O	2.21	0.41
13:CM:106:ASN:O	13:CM:107:ALA:CB	2.69	0.41
1:CA:1501:C:C5	1:CA:1504:G:C5	3.09	0.41
35:DM:132:VAL:HG11	44:DV:81:ARG:NH1	2.36	0.41
23:DA:1647:G:P	23:DA:1647:G:H3'	2.61	0.41
23:BA:1862:G:H2'	23:BA:1863:G:C8	2.54	0.41
23:DA:2013:A:H4'	41:DS:96:ILE:HD12	2.03	0.41
36:BN:104:ARG:CB	36:BN:104:ARG:HH11	2.33	0.41
28:DF:16:ARG:N	28:DF:17:PRO:HD2	2.36	0.41
24:BB:62:C:C2	24:BB:63:G:C8	3.09	0.41
2:AB:137:ARG:O	2:AB:141:GLU:HG2	2.21	0.41
23:DA:1198:U:H2'	23:DA:1199:U:C6	2.56	0.41
23:BA:2836:U:H2'	23:BA:2837:G:H8	1.86	0.41
7:AG:91:VAL:HG12	7:AG:95:ARG:HB3	2.02	0.41
23:BA:2586:C:O5'	23:BA:2586:C:H6	2.04	0.41
23:DA:1229:G:H2'	23:DA:1230:C:H6	1.86	0.41
1:AA:862:C:H2'	1:AA:863:U:O4'	2.20	0.41
45:BW:27:GLU:HB2	45:BW:69:PHE:CD1	2.55	0.41
1:AA:901:A:C5	1:AA:902:G:H1'	2.56	0.41
23:DA:945:A:O2'	23:DA:946:G:H4'	2.21	0.41
23:BA:172:C:H2'	23:BA:173:G:H8	1.86	0.41
1:CA:1448:C:H2'	1:CA:1449:C:C6	2.55	0.41
23:DA:2459:A:C4	23:DA:2460:U:C6	3.09	0.41
1:CA:1352:C:H2'	1:CA:1353:G:C8	2.56	0.41
1:CA:355:C:C4	1:CA:356:A:N7	2.89	0.41
1:CA:809:G:C6	1:CA:810:C:C5	3.09	0.41
23:BA:1831:G:C5	23:BA:1832:C:C5	3.09	0.41
9:CI:33:PHE:HZ	9:CI:43:ALA:O	2.03	0.41
23:BA:952:G:P	35:BM:16:ARG:HH22	2.44	0.41
46:DX:70:VAL:O	46:DX:74:VAL:HG23	2.20	0.41
23:DA:540:G:C4	23:DA:541:C:C5	3.09	0.41
23:BA:2230:G:C6	23:BA:2231:C:C4	3.09	0.41
23:DA:1368:G:C2	23:DA:1369:G:C8	3.09	0.41
23:BA:686:G:N2	23:BA:788:A:H61	2.18	0.41
38:DP:137:LYS:HD2	38:DP:137:LYS:N	2.36	0.41
6:AF:70:ASP:OD1	6:AF:70:ASP:N	2.50	0.41
23:DA:885:C:H6	23:DA:885:C:O5'	2.04	0.41
27:DE:9:ILE:HD13	27:DE:9:ILE:H	1.86	0.41
12:CL:45:LYS:HE2	12:CL:45:LYS:HB3	1.87	0.41
27:DE:24:LEU:HA	27:DE:25:PRO:HD3	1.86	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:874:G:C5	1:CA:875:C:C5	3.09	0.41
1:AA:1095:U:H2'	1:AA:1096:C:C6	2.56	0.41
23:DA:941:A:H4'	34:DL:35:HIS:CD2	2.56	0.40
1:CA:1399:C:C2	1:CA:1401:G:C6	3.09	0.40
30:BH:92:VAL:HG22	30:BH:92:VAL:O	2.20	0.40
43:DU:30:VAL:HG13	43:DU:37:VAL:HG12	2.03	0.40
23:BA:2685:G:N3	23:BA:2725:A:C2	2.89	0.40
1:CA:735:C:O2'	1:CA:736:C:H5'	2.20	0.40
1:AA:738:C:H2'	1:AA:739:C:H6	1.85	0.40
25:BC:197:GLY:O	25:BC:198:ASN:C	2.59	0.40
44:BV:48:PHE:HE2	44:BV:71:VAL:HG21	1.86	0.40
44:BV:97:GLU:O	44:BV:98:MET:HB3	2.20	0.40
25:DC:133:LEU:HB3	25:DC:173:VAL:HG11	2.02	0.40
23:BA:2028:U:H2'	23:BA:2029:G:C8	2.55	0.40
23:BA:1509:A:H4'	23:BA:1510:A:N9	2.36	0.40
23:BA:1174:A:H3'	23:BA:1175:U:C5'	2.46	0.40
34:BL:122:PRO:O	34:BL:123:LEU:HB3	2.21	0.40
23:DA:2343:C:O2'	23:DA:2344:U:H5'	2.21	0.40
36:BN:14:SER:O	36:BN:15:SER:C	2.60	0.40
23:BA:972:G:H3'	23:BA:973:A:H2'	2.04	0.40
33:DK:106:LEU:N	33:DK:106:LEU:HD12	2.36	0.40
23:BA:646:A:H2'	23:BA:647:G:O4'	2.20	0.40
13:AM:3:ARG:HG2	13:AM:9:ILE:CD1	2.51	0.40
25:DC:118:VAL:HG13	25:DC:119:ALA:N	2.37	0.40
46:BX:84:GLY:O	46:BX:85:LEU:C	2.59	0.40
35:DM:60:ARG:HH11	35:DM:60:ARG:HB2	1.86	0.40
4:AD:172:PRO:HD2	4:AD:173:TRP:CZ3	2.56	0.40
4:CD:23:GLY:HA3	4:CD:112:VAL:CG2	2.49	0.40
23:BA:1858:G:HO2'	23:BA:1859:A:H8	1.66	0.40
33:BK:7:TYR:CE1	33:BK:20:MET:HB2	2.56	0.40
9:CI:112:LYS:HA	9:CI:119:ALA:HB2	2.03	0.40
24:DB:73:A:C4	24:DB:104:A:C2	3.09	0.40
34:DL:21:ARG:H	34:DL:21:ARG:HG2	1.73	0.40
23:BA:2777:G:C4'	23:BA:2778:A:H5'	2.51	0.40
25:DC:79:VAL:HG11	25:DC:111:LEU:CD1	2.51	0.40
29:DG:104:GLU:HA	29:DG:113:VAL:O	2.22	0.40
36:BN:11:ASN:CG	36:BN:12:ARG:H	2.21	0.40
23:DA:784:A:H5'	23:DA:785:G:OP1	2.20	0.40
38:BP:19:LEU:HA	38:BP:20:PRO:HD3	1.73	0.40
13:AM:115:LYS:HE3	13:AM:115:LYS:HB2	1.86	0.40
1:CA:17:U:N3	1:CA:18:C:C4	2.89	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BA:276:A:C3'	23:BA:277:C:H5''	2.51	0.40
1:CA:16:A:N1	1:CA:919:A:H2	2.19	0.40
23:BA:2749:A:H1'	29:BG:63:SER:OG	2.21	0.40
13:CM:91:ARG:CD	19:CS:81:ARG:HH22	2.35	0.40
23:DA:524:U:O2'	23:DA:554:U:H4'	2.21	0.40
37:DO:38:GLN:HB3	37:DO:47:THR:HG21	2.03	0.40
41:BS:82:LEU:HD23	41:BS:84:ARG:NH2	2.36	0.40
30:DH:66:GLU:O	30:DH:70:GLU:HG2	2.20	0.40
13:CM:86:CYS:SG	13:CM:88:ARG:HB2	2.61	0.40
36:BN:28:LEU:HD23	36:BN:34:ILE:HG12	2.02	0.40
23:DA:2626:C:H2'	23:DA:2627:G:O4'	2.21	0.40
23:BA:1936:A:H5''	23:BA:1936:A:N3	2.36	0.40
8:CH:17:THR:C	8:CH:78:GLN:HE22	2.24	0.40
27:DE:47:GLY:O	27:DE:94:PRO:HB3	2.21	0.40
40:DR:89:GLN:HA	40:DR:90:PRO:HD3	1.80	0.40
35:BM:34:LEU:HD12	35:BM:130:LYS:O	2.21	0.40
1:AA:658:G:C6	1:AA:659:U:C4	3.09	0.40
23:BA:2018:G:C6	23:BA:2019:A:C6	3.09	0.40
23:BA:1726:G:C2	23:BA:1735:U:O2	2.74	0.40
13:CM:56:LEU:O	13:CM:56:LEU:HD13	2.21	0.40
27:DE:39:TRP:CD1	27:DE:101:LEU:HB2	2.57	0.40
1:CA:711:G:O2'	1:CA:712:A:H5'	2.21	0.40
23:DA:2642:G:O2'	23:DA:2643:G:H5'	2.21	0.40
23:BA:1198:U:H2'	23:BA:1199:U:C6	2.55	0.40
11:CK:34:ASP:HB2	11:CK:35:PRO:HD2	2.03	0.40
1:AA:874:G:C5	1:AA:875:C:C5	3.09	0.40
23:DA:2762:G:O2'	23:DA:2763:G:H5'	2.21	0.40
14:AN:43:CYS:O	14:AN:47:LEU:HG	2.21	0.40
23:BA:300:A:OP1	43:BU:84:ARG:NH2	2.54	0.40
1:CA:667:G:H4'	15:CO:51:HIS:CE1	2.56	0.40
13:AM:58:GLU:OE1	13:AM:62:ASN:HB2	2.21	0.40
23:DA:386:G:H4'	23:DA:387:U:OP2	2.21	0.40
23:BA:1502:C:O5'	23:BA:1502:C:H6	2.04	0.40
10:CJ:22:LYS:HD2	10:CJ:22:LYS:O	2.21	0.40
23:DA:1532:C:O5'	23:DA:1532:C:H6	2.04	0.40
40:DR:6:LYS:HG3	40:DR:6:LYS:O	2.21	0.40
23:DA:2685:G:N3	23:DA:2725:A:C2	2.89	0.40
23:BA:2422:A:C5	23:BA:2424:C:C4	3.09	0.40
35:BM:22:LYS:HA	35:BM:22:LYS:CE	2.35	0.40
23:BA:827:U:H2'	23:BA:2068:U:C2	2.56	0.40
43:DU:97:ARG:HD3	43:DU:98:VAL:HG12	2.04	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:735:C:O2'	1:AA:736:C:H5'	2.21	0.40
35:DM:81:VAL:O	35:DM:82:ARG:HD3	2.21	0.40
46:DX:10:LYS:O	46:DX:11:ARG:CG	2.67	0.40
29:DG:86:GLU:HB3	29:DG:132:ARG:NH1	2.37	0.40
35:BM:141:GLN:O	44:BV:53:ILE:HG22	2.20	0.40
1:CA:105:G:C4	1:CA:106:C:C5	3.09	0.40
45:BW:32:ARG:CA	45:BW:35:ASN:HD21	2.34	0.40
25:DC:131:LEU:HD12	25:DC:136:ILE:HG12	2.02	0.40
1:AA:691:G:H3'	11:AK:26:ASN:ND2	2.35	0.40
36:BN:4:LEU:C	36:BN:6:SER:N	2.72	0.40
2:AB:187:LEU:HD11	2:AB:205:ASP:HB3	2.04	0.40
53:B5:13:ARG:HG3	53:B5:14:VAL:HG23	2.02	0.40
23:BA:2029:G:H2'	23:BA:2031:A:OP1	2.22	0.40
23:BA:603:A:C2	23:BA:655:A:N3	2.89	0.40
28:BF:53:LEU:HD13	28:BF:88:ILE:HG12	2.03	0.40
2:AB:27:LYS:O	2:AB:30:ARG:HG2	2.21	0.40
36:BN:14:SER:HA	36:BN:17:ARG:HG2	2.03	0.40
26:BD:5:LEU:HD22	26:BD:197:ILE:HG22	2.02	0.40
4:AD:49:ARG:CZ	4:AD:50:ARG:HB2	2.51	0.40
30:BH:114:LEU:HA	30:BH:130:TYR:HD1	1.86	0.40
30:BH:128:LEU:HG	30:BH:142:VAL:CG2	2.50	0.40
44:DV:144:LEU:HD22	44:DV:144:LEU:N	2.36	0.40
23:DA:388:G:C4	23:DA:390:A:C6	3.10	0.40
1:CA:1226:C:C4	13:CM:104:ARG:HB2	2.57	0.40
11:AK:109:VAL:HG12	11:AK:110:ASP:N	2.36	0.40
1:AA:542:G:H5'	4:AD:41:GLY:HA2	2.03	0.40
34:DL:126:VAL:HG22	34:DL:145:PRO:HB2	2.02	0.40
23:BA:1276:A:H5''	23:BA:1276:A:H8	1.86	0.40
5:AE:92:LYS:HA	5:AE:93:PRO:HD2	1.89	0.40
1:CA:187:C:O2	1:CA:191(A):G:C6	2.74	0.40
47:BY:24:LEU:HD23	47:BY:24:LEU:O	2.21	0.40
23:BA:265:A:C8	23:BA:266:G:H1'	2.55	0.40
23:BA:534:U:H1'	39:BQ:49:HIS:CD2	2.56	0.40
46:BX:23:LYS:HG3	46:BX:23:LYS:O	2.20	0.40
44:DV:161:VAL:HG12	44:DV:162:GLU:N	2.35	0.40
1:CA:815:A:H4'	1:CA:817:C:C5	2.56	0.40
1:AA:555:C:H2'	1:AA:556:C:H6	1.84	0.40
51:B3:23:THR:O	51:B3:24:GLU:HG2	2.22	0.40
23:DA:1639:U:C2'	23:DA:1640:C:H5''	2.51	0.40
23:DA:2293:C:H4'	37:DO:93:LYS:HZ2	1.87	0.40
44:DV:118:GLN:HB2	44:DV:173:ALA:C	2.42	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
13:AM:91:ARG:CD	19:AS:81:ARG:HH22	2.35	0.40
27:BE:40:GLN:O	27:BE:43:LYS:HG2	2.21	0.40
25:BC:174:ILE:CD1	25:BC:174:ILE:N	2.84	0.40
23:BA:69:C:H2'	23:BA:70:G:H8	1.84	0.40
28:BF:16:ARG:N	28:BF:17:PRO:HD2	2.36	0.40
32:BJ:159:GLU:HG2	32:BJ:160:LYS:N	2.36	0.40
23:BA:1271:G:O3'	23:BA:1272:A:H4'	2.21	0.40
23:DA:1289:C:H2'	23:DA:1290:C:C6	2.55	0.40
2:CB:137:ARG:O	2:CB:141:GLU:HG2	2.21	0.40
23:DA:451:C:N4	23:DA:453:C:H3'	2.36	0.40
23:DA:1270:C:H5''	23:DA:1271:G:C5'	2.52	0.40
1:CA:901:A:C5	1:CA:902:G:H1'	2.55	0.40
1:AA:838:G:N2	1:AA:849:C:N3	2.70	0.40
1:AA:1010:G:H2'	1:AA:1011:G:H8	1.87	0.40
23:DA:570:G:H2'	23:DA:2030:A:C5	2.55	0.40
23:BA:1203:G:C6	23:BA:1204:A:N6	2.89	0.40
1:AA:552:U:H4'	12:AL:85:ARG:HG2	2.04	0.40
35:DM:67:ARG:HD2	35:DM:105:GLU:OE2	2.21	0.40
1:CA:1437:C:H2'	1:CA:1438:G:C8	2.56	0.40
1:AA:987:G:H2'	1:AA:988:G:C8	2.56	0.40
7:AG:80:VAL:C	7:AG:82:GLY:H	2.24	0.40
1:CA:1269:A:H5'	21:CU:19:GLY:HA2	2.04	0.40
1:AA:936:C:H2'	1:AA:937:A:O4'	2.22	0.40
23:BA:1918:A:O2'	23:BA:1920:C:N4	2.54	0.40
32:BJ:132:LYS:HG2	32:BJ:132:LYS:H	1.61	0.40
42:DT:41:ASN:N	42:DT:41:ASN:HD22	2.19	0.40
46:DX:82:LEU:N	46:DX:82:LEU:HD12	2.36	0.40
4:AD:200:GLU:H	4:AD:200:GLU:CD	2.24	0.40
4:AD:209:ARG:HA	4:AD:209:ARG:HD2	1.90	0.40
1:AA:1437:C:H2'	1:AA:1438:G:C8	2.56	0.40
35:BM:67:ARG:HD2	35:BM:105:GLU:OE2	2.21	0.40
1:CA:1046:A:H3'	1:CA:1047:G:H8	1.86	0.40
11:CK:23:ALA:HB1	11:CK:88:GLY:HA3	2.03	0.40
23:BA:252:G:OP2	34:BL:50:ARG:NH2	2.49	0.40
34:BL:39:LYS:CD	34:BL:40:SER:H	2.35	0.40
39:DQ:83:LEU:N	39:DQ:83:LEU:HD12	2.36	0.40
25:BC:236:GLY:O	25:BC:238:GLY:N	2.54	0.40
23:DA:2246:G:H2'	23:DA:2247:A:H8	1.87	0.40
1:AA:954:G:H2'	1:AA:955:U:C6	2.56	0.40
34:BL:62:LEU:HD22	34:BL:62:LEU:O	2.22	0.40
23:BA:2416:C:OP1	34:BL:64:LYS:O	2.40	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
46:BX:46:LEU:HB2	46:BX:63:ALA:HA	2.04	0.40
13:AM:99:ARG:HB2	13:AM:101:GLN:NE2	2.26	0.40
1:AA:692:U:C5	11:AK:26:ASN:ND2	2.87	0.40
26:DD:184:VAL:HG12	26:DD:185:LYS:H	1.86	0.40
23:BA:1493:C:O2	23:BA:1493:C:H2'	2.21	0.40
23:DA:1509:A:H4'	23:DA:1510:A:N9	2.36	0.40
12:CL:74:HIS:HD2	12:CL:76:LEU:HB2	1.85	0.40
2:CB:80:ILE:HD11	2:CB:208:ILE:CG2	2.50	0.40
24:DB:70:C:H2'	24:DB:71:C:C6	2.52	0.40
36:BN:10:LEU:HD12	36:BN:10:LEU:N	2.37	0.40
11:CK:20:TYR:C	11:CK:21:ILE:HD12	2.41	0.40
46:BX:41:ARG:HD3	46:BX:43:TYR:CE2	2.57	0.40
34:BL:17:LYS:C	34:BL:19:VAL:H	2.25	0.40
19:CS:6:LYS:HD2	19:CS:6:LYS:N	2.34	0.40
23:DA:643:A:OP1	51:D3:42:TRP:NE1	2.54	0.40
44:BV:144:LEU:HD22	44:BV:144:LEU:N	2.37	0.40
28:DF:7:LEU:HA	28:DF:10:LYS:HD2	2.03	0.40
4:AD:188:LEU:HD12	4:AD:188:LEU:H	1.86	0.40
41:DS:18:ARG:NH1	41:DS:76:VAL:HG13	2.36	0.40
23:BA:1858:G:H1'	23:BA:1884:A:H62	1.86	0.40
1:CA:1079:G:C6	1:CA:1080:A:N6	2.89	0.40
1:AA:438:G:C4'	1:AA:439:A:OP1	2.67	0.40
23:BA:480:A:H2'	23:BA:480:A:N3	2.36	0.40
46:BX:73:LEU:HD21	46:BX:94:LEU:CG	2.49	0.40
19:AS:63:THR:HG22	19:AS:66:MET:CE	2.51	0.40
29:DG:103:LEU:HG	29:DG:105:LEU:CD1	2.51	0.40
25:BC:107:ALA:HA	25:BC:108:PRO:HD3	1.78	0.40
7:CG:150:ALA:HA	11:CK:59:TYR:CD2	2.56	0.40
32:DJ:83:ILE:HG22	32:DJ:84:ARG:O	2.21	0.40
28:BF:114:ILE:HB	28:BF:117:PHE:HB2	2.02	0.40
23:DA:656:G:H8	23:DA:656:G:O5'	2.05	0.40
23:BA:839:U:H2'	23:BA:840:C:H6	1.81	0.40
23:BA:775:G:C2	23:BA:777:A:N6	2.89	0.40
23:DA:830:G:C4	23:DA:2448:A:C6	3.09	0.40
2:CB:102:LEU:HD12	2:CB:102:LEU:N	2.37	0.40
1:AA:57:G:C6	1:AA:58:C:C4	3.09	0.40
9:CI:53:VAL:HG11	9:CI:85:LEU:HD22	2.02	0.40
1:CA:235:C:H1'	17:CQ:61:GLU:OE1	2.21	0.40
44:DV:127:LYS:HB3	44:DV:162:GLU:CG	2.51	0.40
1:AA:161:A:H2'	1:AA:162:A:H8	1.85	0.40
13:AM:8:GLU:OE1	13:AM:22:ILE:HG23	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
13:CM:19:LEU:HA	13:CM:22:ILE:HG12	2.02	0.40
16:CP:49:LEU:HD12	16:CP:50:LYS:N	2.36	0.40
35:DM:77:LYS:HA	35:DM:78:PRO:HD3	1.73	0.40
1:CA:1053:G:C3'	1:CA:1054:C:H5'	2.51	0.40
23:BA:1344:G:H5'	23:BA:1384:A:C6	2.56	0.40
8:CH:82:HIS:HD2	8:CH:138:TRP:HE1	1.69	0.40
1:AA:562:C:H5'	1:AA:562:C:H6	1.85	0.40
26:BD:76:ARG:HG2	26:BD:77:ILE:HG13	2.03	0.40
23:BA:1599:C:H2'	23:BA:1600:C:C6	2.56	0.40
23:BA:1488:G:C5	23:BA:1489:U:C5	3.10	0.40
25:DC:175:LEU:HD23	25:DC:175:LEU:HA	1.95	0.40
1:AA:900:A:H2'	1:AA:901:A:C8	2.56	0.40
46:BX:53:VAL:HG22	46:BX:74:VAL:HG13	2.03	0.40
13:CM:37:THR:OG1	13:CM:56:LEU:HD23	2.22	0.40
1:CA:605:U:H2'	1:CA:606:G:O4'	2.22	0.40
6:CF:9:VAL:HG13	6:CF:59:TYR:O	2.21	0.40
1:CA:69:G:H2'	1:CA:73:G:H8	1.86	0.40
23:DA:327:G:C2	23:DA:328:U:C2	3.10	0.40
23:DA:2386:C:H4'	45:DW:55:ARG:O	2.21	0.40
23:BA:386:G:H4'	23:BA:387:U:OP2	2.22	0.40
23:BA:2404:C:H2'	23:BA:2405:G:O4'	2.22	0.40
20:CT:36:LEU:HB3	20:CT:59:ALA:HB2	2.03	0.40
29:DG:117:PRO:HA	29:DG:118:PRO:HD2	1.93	0.40
23:BA:1221:C:H2'	23:BA:122(A):C:C6	2.56	0.40
10:AJ:22:LYS:HD2	10:AJ:22:LYS:O	2.21	0.40
1:AA:604:G:C6	1:AA:605:U:C4	3.09	0.40
23:BA:2574:G:H8	23:BA:2574:G:O5'	2.04	0.40
23:BA:1728:G:H8	23:BA:1728:G:O5'	2.05	0.40
38:BP:137:LYS:N	38:BP:137:LYS:HD2	2.36	0.40
16:CP:12:LYS:C	16:CP:14:ASN:H	2.24	0.40
1:AA:712:A:O2'	1:AA:713:G:H5'	2.21	0.40
7:AG:54:THR:C	7:AG:56:GLN:H	2.25	0.40
23:BA:587:C:H4'	23:BA:588:U:C6	2.57	0.40
34:BL:47:ASP:HB3	34:BL:48:PRO:HA	2.02	0.40
37:BO:82:ILE:HG22	37:BO:83:LYS:N	2.35	0.40
51:D3:25:LYS:HD3	53:D5:34:TRP:CH2	2.56	0.40
47:DY:1:MET:SD	47:DY:5:GLU:OE2	2.80	0.40
23:BA:97:C:H5''	47:BY:2:LYS:HZ1	1.86	0.40
23:DA:270(I):C:O2'	23:DA:270(J):G:H5'	2.22	0.40
44:DV:71:VAL:HG11	44:DV:74:VAL:HG23	2.02	0.40
11:CK:44:SER:H	11:CK:47:VAL:HB	1.87	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:DL:57:THR:C	34:DL:59:LEU:N	2.75	0.40
30:BH:133:HIS:HA	30:BH:134:PRO:HD3	1.92	0.40
23:BA:1310:G:H2'	23:BA:1311:G:H5''	2.03	0.40
23:DA:1408:C:N3	23:DA:1595:G:C2	2.89	0.40
23:BA:141(A):A:C8	23:BA:1408:C:H1'	2.57	0.40
23:DA:1211:U:H4'	23:DA:1212:G:OP2	2.22	0.40
33:DK:71:ARG:NH2	33:DK:77:ILE:HG21	2.37	0.40
30:DH:83:ALA:N	30:DH:89:TYR:HD1	2.20	0.40
13:CM:3:ARG:HG2	13:CM:9:ILE:CD1	2.52	0.40
1:AA:1320:C:C2	19:AS:72:GLY:HA3	2.56	0.40
23:DA:2484:G:H2'	23:DA:2485:G:H8	1.86	0.40
28:BF:171:ALA:O	28:BF:175:LEU:HG	2.21	0.40
22:CV:6188:C:HO2'	22:CV:6189:A:P	2.45	0.40
30:DH:130:TYR:HD2	30:DH:132:PRO:HG3	1.83	0.40
1:AA:1296:C:OP1	13:AM:44:ARG:NH2	2.51	0.40
40:BR:72:VAL:O	40:BR:72:VAL:HG23	2.22	0.40
7:AG:15:ASP:HB2	7:AG:20:ASP:O	2.21	0.40
23:DA:443:A:H1'	23:DA:1201:C:O4'	2.21	0.40
1:CA:1079:G:O3'	5:CE:14:ARG:NH2	2.54	0.40
32:DJ:117:HIS:HA	32:DJ:118:PRO:HD2	1.98	0.40
23:BA:270(H):C:H2'	23:BA:270(I):C:C6	2.56	0.40
26:BD:85:ASN:HA	26:BD:86:PRO:HD3	1.78	0.40
46:DX:35:THR:HB	46:DX:36:GLY:H	1.54	0.40
1:CA:506:G:C6	1:CA:507:C:N4	2.89	0.40
23:BA:2477:C:HO2'	23:BA:2478:A:P	2.44	0.40
23:BA:2565:A:H5''	23:BA:2566:A:OP2	2.22	0.40
1:CA:892:A:C6	1:CA:893:C:C4	3.09	0.40
1:AA:1501:C:C5	1:AA:1504:G:C5	3.09	0.40
41:BS:50:VAL:HG11	41:BS:103:ILE:HG21	2.04	0.40
24:BB:37:C:C5	24:BB:38:C:C5	3.09	0.40
35:BM:78:PRO:C	35:BM:79:LEU:HD12	2.42	0.40
23:DA:2230:G:H1'	46:DX:45:ASN:CB	2.51	0.40
38:DP:3:ARG:NH1	38:DP:6:LEU:HD23	2.37	0.40
23:BA:118:A:H1'	23:BA:178:G:O4'	2.22	0.40
23:BA:875:G:N2	23:BA:903:C:C2	2.90	0.40
28:BF:133:LEU:HD23	28:BF:133:LEU:N	2.36	0.40
28:BF:39:ILE:HG23	28:BF:157:ILE:HG22	2.04	0.40
1:AA:620:C:C2	4:AD:135:LEU:HG	2.56	0.40
27:DE:72:ARG:O	27:DE:73:ALA:O	2.39	0.40
49:B1:40:ILE:HD12	49:B1:40:ILE:N	2.36	0.40
29:BG:83:TYR:CZ	29:BG:138:LYS:HG3	2.56	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BA:2019:A:H5''	39:BQ:27:LEU:HD12	2.04	0.40
1:CA:987:G:H2'	1:CA:988:G:C8	2.57	0.40
23:DA:991:C:C5	23:DA:1185:C:C4	3.09	0.40
23:BA:1843:C:O2'	23:BA:1844:C:H5'	2.22	0.40
34:DL:106:LEU:HA	34:DL:106:LEU:HD22	1.86	0.40
43:BU:98:VAL:HG22	43:BU:98:VAL:O	2.21	0.40
27:DE:108:LYS:HD3	27:DE:108:LYS:HA	1.94	0.40
27:DE:64:ILE:HD12	27:DE:64:ILE:HA	1.81	0.40
35:DM:80:GLU:HA	35:DM:80:GLU:OE2	2.21	0.40
30:DH:136:VAL:N	30:DH:137:PRO:HD3	2.37	0.40
23:BA:1783:A:C2	23:BA:2587:A:C4	3.10	0.40
23:DA:2338:G:C2	23:DA:2339:G:C8	3.10	0.40
23:DA:2065:C:H2'	23:DA:2066:C:C6	2.56	0.40
1:CA:1051:C:C4	1:CA:1052:U:C4	3.09	0.40
23:DA:1401:G:H2'	23:DA:1402:C:C6	2.57	0.40
23:BA:2057:A:N6	23:BA:2058:A:C6	2.89	0.40
39:DQ:91:ASP:OD2	39:DQ:96:ALA:HB2	2.22	0.40
23:DA:782:A:H5'	23:DA:783:A:C2	2.57	0.40
32:DJ:157:ARG:HG2	32:DJ:157:ARG:O	2.21	0.40
34:DL:114:ILE:HD11	34:DL:130:PHE:CG	2.54	0.40
34:DL:64:LYS:O	34:DL:65:ARG:C	2.59	0.40
38:DP:26:ASP:HB3	38:DP:92:GLY:H	1.86	0.40
1:CA:942:G:H2'	1:CA:943:U:H6	1.86	0.40
3:AC:182:ILE:HG23	3:AC:202:ILE:C	2.41	0.40
25:BC:35:LYS:O	25:BC:64:ILE:HD12	2.21	0.40
34:DL:122:PRO:O	34:DL:123:LEU:HB3	2.21	0.40
30:BH:88:ILE:HG22	30:BH:89:TYR:N	2.35	0.40
23:DA:2287:A:C5	23:DA:2289:G:C5	3.10	0.40
28:DF:87:PRO:HB2	28:DF:88:ILE:H	1.65	0.40
30:BH:114:LEU:HA	30:BH:130:TYR:CD1	2.57	0.40
23:BA:468:G:N7	52:B4:39:ARG:NH2	2.68	0.40
27:BE:155:LEU:HA	27:BE:174:VAL:HG23	2.03	0.40
37:BO:26:LEU:HB3	37:BO:87:PHE:HA	2.03	0.40
37:BO:87:PHE:HE2	37:BO:89:ARG:HA	1.87	0.40
37:DO:26:LEU:HB3	37:DO:87:PHE:HA	2.03	0.40
4:CD:108:LEU:HD12	4:CD:108:LEU:HA	1.94	0.40
11:CK:26:ASN:O	11:CK:27:ASN:HB2	2.21	0.40
36:DN:38:VAL:HG22	36:DN:112:ALA:HB2	2.04	0.40
1:CA:272:C:H2'	1:CA:273:A:H8	1.86	0.40
23:DA:7:G:H2'	23:DA:8:A:O4'	2.21	0.40
19:AS:45:VAL:HA	19:AS:62:ILE:HG23	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DA:444:C:H2'	23:DA:445:C:H6	1.86	0.40
32:DJ:64:ASP:N	32:DJ:64:ASP:OD1	2.54	0.40
11:CK:59:TYR:O	11:CK:62:GLN:HB3	2.22	0.40
41:DS:29:LEU:HD22	41:DS:69:LEU:CD1	2.49	0.40
1:AA:1316:G:H22	1:AA:1319:A:H5''	1.86	0.40
1:AA:300:A:C8	1:AA:300:A:H3'	2.57	0.40
9:AI:99:LEU:HD12	9:AI:101:PHE:CE2	2.56	0.40
19:CS:45:VAL:HA	19:CS:62:ILE:HG23	2.02	0.40
23:DA:758:C:O2	23:DA:1981:A:H2	2.04	0.40
38:BP:128:GLU:O	38:BP:132:LYS:HG3	2.20	0.40
38:BP:92:GLY:HA2	38:BP:117:ASP:H	1.87	0.40
23:DA:319:C:C2	23:DA:320:A:C8	3.08	0.40
1:AA:663:A:H2'	1:AA:664:G:O4'	2.22	0.40
23:DA:2565:A:H5''	23:DA:2566:A:OP2	2.22	0.40
1:CA:1234:C:O2'	1:CA:1235:U:H5'	2.22	0.40
1:AA:1228:C:H5'	13:AM:115:LYS:O	2.21	0.40
9:AI:92:TYR:O	9:AI:96:LEU:HB2	2.22	0.40
35:DM:27:VAL:H	44:DV:81:ARG:NH2	2.20	0.40
35:BM:27:VAL:H	44:BV:81:ARG:NH2	2.19	0.40
23:BA:1647:G:H3'	23:BA:1647:G:P	2.62	0.40
23:BA:319:C:C2	23:BA:320:A:C8	3.10	0.40
44:DV:145:GLU:HG3	44:DV:146:ILE:N	2.37	0.40
24:DB:48:A:H2'	24:DB:49:C:C6	2.57	0.40
2:AB:130:ARG:HA	2:AB:131:PRO:HD2	1.95	0.40
1:CA:928:G:C2	1:CA:1390:U:O2	2.74	0.40
26:BD:14:ILE:HD12	26:BD:14:ILE:O	2.22	0.40
23:DA:380:U:H2'	23:DA:381:G:H8	1.87	0.40
23:BA:2846:G:C5	23:BA:2847:U:C5	3.09	0.40
23:DA:1710:C:O2'	23:DA:1711:C:H5'	2.22	0.40
23:DA:1750:G:H2'	23:DA:1751:C:H6	1.87	0.40
18:CR:45:SER:HB3	18:CR:51:LEU:HD21	2.03	0.40
44:BV:28:MET:HA	44:BV:88:PHE:HB2	2.03	0.40
23:BA:2764:A:N6	23:BA:2766:G:C2	2.90	0.40
1:CA:179:A:C4	1:CA:180:U:C5	3.09	0.40
23:BA:489:G:C5	23:BA:1284:A:C2	3.09	0.40
2:CB:16:HIS:CD2	2:CB:210:SER:HA	2.57	0.40
1:AA:1250:A:H2	1:AA:1353:G:H21	1.69	0.40
6:AF:95:GLU:HA	6:AF:96:PRO:HD3	1.96	0.40
2:AB:46:LYS:HD2	2:AB:49:GLU:OE1	2.22	0.40
1:AA:1046:A:H3'	1:AA:1047:G:H8	1.87	0.40
23:DA:2489:G:C6	23:DA:2490:G:N1	2.89	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
29:DG:125:VAL:HG22	29:DG:131:VAL:HG22	2.03	0.40
1:CA:1259:C:O5'	1:CA:1259:C:H6	2.03	0.40
36:BN:99:LYS:HE3	36:BN:99:LYS:HB3	1.87	0.40
1:AA:608:A:C4	1:AA:609:A:C8	3.09	0.40
8:AH:20:TYR:HA	8:AH:65:TYR:CE2	2.57	0.40
33:BK:90:GLN:O	33:BK:91:LEU:HB2	2.21	0.40
11:AK:94:ALA:O	11:AK:98:LEU:HG	2.21	0.40
23:DA:1961:C:O2'	23:DA:1962:C:H5'	2.22	0.40

There are no symmetry-related clashes.

## 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	AB	232/234 (99%)	194 (84%)	32 (14%)	6 (3%)	7	45
2	CB	232/234 (99%)	194 (84%)	32 (14%)	6 (3%)	7	45
3	AC	204/206 (99%)	153 (75%)	38 (19%)	13 (6%)	2	20
3	CC	204/206 (99%)	154 (76%)	37 (18%)	13 (6%)	2	20
4	AD	206/208 (99%)	172 (84%)	26 (13%)	8 (4%)	4	34
4	CD	206/208 (99%)	172 (84%)	25 (12%)	9 (4%)	3	30
5	AE	149/151 (99%)	123 (83%)	22 (15%)	4 (3%)	6	44
5	CE	149/151 (99%)	126 (85%)	19 (13%)	4 (3%)	6	44
6	AF	99/101 (98%)	90 (91%)	8 (8%)	1 (1%)	19	66
6	CF	99/101 (98%)	90 (91%)	8 (8%)	1 (1%)	19	66
7	AG	153/155 (99%)	132 (86%)	18 (12%)	3 (2%)	9	51
7	CG	153/155 (99%)	133 (87%)	17 (11%)	3 (2%)	9	51
8	AH	136/138 (99%)	119 (88%)	17 (12%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
8	CH	136/138 (99%)	119 (88%)	17 (12%)	0	100	100
9	AI	125/127 (98%)	102 (82%)	20 (16%)	3 (2%)	7	47
9	CI	125/127 (98%)	101 (81%)	21 (17%)	3 (2%)	7	47
10	AJ	96/98 (98%)	78 (81%)	12 (12%)	6 (6%)	2	21
10	CJ	96/98 (98%)	78 (81%)	12 (12%)	6 (6%)	2	21
11	AK	117/119 (98%)	97 (83%)	16 (14%)	4 (3%)	5	39
11	CK	117/119 (98%)	97 (83%)	16 (14%)	4 (3%)	5	39
12	AL	122/124 (98%)	94 (77%)	22 (18%)	6 (5%)	3	27
12	CL	122/124 (98%)	94 (77%)	22 (18%)	6 (5%)	3	27
13	AM	114/116 (98%)	94 (82%)	16 (14%)	4 (4%)	4	38
13	CM	114/116 (98%)	94 (82%)	16 (14%)	4 (4%)	4	38
14	AN	58/60 (97%)	50 (86%)	5 (9%)	3 (5%)	2	25
14	CN	58/60 (97%)	50 (86%)	5 (9%)	3 (5%)	2	25
15	AO	86/88 (98%)	77 (90%)	7 (8%)	2 (2%)	8	48
15	CO	86/88 (98%)	77 (90%)	7 (8%)	2 (2%)	8	48
16	AP	81/83 (98%)	65 (80%)	14 (17%)	2 (2%)	7	46
16	CP	81/83 (98%)	65 (80%)	13 (16%)	3 (4%)	4	36
17	AQ	97/99 (98%)	81 (84%)	15 (16%)	1 (1%)	19	66
17	CQ	97/99 (98%)	81 (84%)	15 (16%)	1 (1%)	19	66
18	AR	68/70 (97%)	53 (78%)	13 (19%)	2 (3%)	6	42
18	CR	68/70 (97%)	53 (78%)	12 (18%)	3 (4%)	3	30
19	AS	76/78 (97%)	57 (75%)	14 (18%)	5 (7%)	1	19
19	CS	76/78 (97%)	57 (75%)	14 (18%)	5 (7%)	1	19
20	AT	97/99 (98%)	82 (84%)	12 (12%)	3 (3%)	5	41
20	CT	97/99 (98%)	82 (84%)	12 (12%)	3 (3%)	5	41
21	AU	22/24 (92%)	17 (77%)	4 (18%)	1 (4%)	3	30
21	CU	22/24 (92%)	17 (77%)	4 (18%)	1 (4%)	3	30
25	BC	269/271 (99%)	220 (82%)	31 (12%)	18 (7%)	1	19
25	DC	269/271 (99%)	218 (81%)	33 (12%)	18 (7%)	1	19
26	BD	202/204 (99%)	168 (83%)	26 (13%)	8 (4%)	4	33
26	DD	202/204 (99%)	167 (83%)	29 (14%)	6 (3%)	5	42

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
27	BE	200/202 (99%)	165 (82%)	28 (14%)	7 (4%)	4	38
27	DE	200/202 (99%)	165 (82%)	28 (14%)	7 (4%)	4	38
28	BF	179/181 (99%)	134 (75%)	37 (21%)	8 (4%)	3	30
28	DF	179/181 (99%)	133 (74%)	37 (21%)	9 (5%)	3	27
29	BG	157/159 (99%)	126 (80%)	27 (17%)	4 (2%)	7	46
29	DG	157/159 (99%)	125 (80%)	28 (18%)	4 (2%)	7	46
30	BH	143/145 (99%)	109 (76%)	28 (20%)	6 (4%)	3	32
30	DH	143/145 (99%)	109 (76%)	28 (20%)	6 (4%)	3	32
31	BI	28/65 (43%)	27 (96%)	1 (4%)	0	100	100
31	DI	28/65 (43%)	27 (96%)	1 (4%)	0	100	100
32	BJ	135/137 (98%)	108 (80%)	19 (14%)	8 (6%)	2	22
32	DJ	135/137 (98%)	108 (80%)	19 (14%)	8 (6%)	2	22
33	BK	120/122 (98%)	105 (88%)	8 (7%)	7 (6%)	2	23
33	DK	120/122 (98%)	107 (89%)	6 (5%)	7 (6%)	2	23
34	BL	144/146 (99%)	94 (65%)	35 (24%)	15 (10%)	1	8
34	DL	144/146 (99%)	92 (64%)	37 (26%)	15 (10%)	1	8
35	BM	134/136 (98%)	98 (73%)	22 (16%)	14 (10%)	1	8
35	DM	134/136 (98%)	97 (72%)	24 (18%)	13 (10%)	1	10
36	BN	115/117 (98%)	97 (84%)	13 (11%)	5 (4%)	3	31
36	DN	115/117 (98%)	97 (84%)	14 (12%)	4 (4%)	4	38
37	BO	96/98 (98%)	65 (68%)	18 (19%)	13 (14%)	0	4
37	DO	96/98 (98%)	65 (68%)	18 (19%)	13 (14%)	0	4
38	BP	135/137 (98%)	99 (73%)	30 (22%)	6 (4%)	3	30
38	DP	135/137 (98%)	99 (73%)	30 (22%)	6 (4%)	3	30
39	BQ	114/116 (98%)	99 (87%)	13 (11%)	2 (2%)	11	53
39	DQ	114/116 (98%)	99 (87%)	13 (11%)	2 (2%)	11	53
40	BR	99/101 (98%)	71 (72%)	19 (19%)	9 (9%)	1	11
40	DR	99/101 (98%)	71 (72%)	19 (19%)	9 (9%)	1	11
41	BS	110/112 (98%)	94 (86%)	14 (13%)	2 (2%)	11	53
41	DS	110/112 (98%)	94 (86%)	14 (13%)	2 (2%)	11	53
42	BT	90/92 (98%)	82 (91%)	7 (8%)	1 (1%)	17	63

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
42	DT	90/92 (98%)	82 (91%)	7 (8%)	1 (1%)	17	63
43	BU	98/100 (98%)	65 (66%)	22 (22%)	11 (11%)	0	7
43	DU	98/100 (98%)	63 (64%)	23 (24%)	12 (12%)	0	6
44	BV	186/188 (99%)	140 (75%)	36 (19%)	10 (5%)	2	25
44	DV	186/188 (99%)	140 (75%)	36 (19%)	10 (5%)	2	25
45	BW	74/76 (97%)	59 (80%)	12 (16%)	3 (4%)	3	33
45	DW	74/76 (97%)	57 (77%)	14 (19%)	3 (4%)	3	33
46	BX	86/88 (98%)	57 (66%)	20 (23%)	9 (10%)	1	8
46	DX	86/88 (98%)	57 (66%)	20 (23%)	9 (10%)	1	8
47	BY	60/62 (97%)	48 (80%)	9 (15%)	3 (5%)	3	27
47	DY	60/62 (97%)	49 (82%)	8 (13%)	3 (5%)	3	27
48	BZ	57/59 (97%)	51 (90%)	5 (9%)	1 (2%)	11	53
48	DZ	57/59 (97%)	51 (90%)	5 (9%)	1 (2%)	11	53
49	B1	28/30 (93%)	15 (54%)	10 (36%)	3 (11%)	0	8
49	D1	28/30 (93%)	15 (54%)	10 (36%)	3 (11%)	0	8
50	B2	50/52 (96%)	39 (78%)	8 (16%)	3 (6%)	2	21
50	D2	50/52 (96%)	40 (80%)	6 (12%)	4 (8%)	1	13
51	B3	42/44 (96%)	35 (83%)	2 (5%)	5 (12%)	0	6
51	D3	42/44 (96%)	35 (83%)	2 (5%)	5 (12%)	0	6
52	B4	46/48 (96%)	42 (91%)	4 (9%)	0	100	100
52	D4	46/48 (96%)	41 (89%)	5 (11%)	0	100	100
53	B5	61/63 (97%)	43 (70%)	13 (21%)	5 (8%)	1	13
53	D5	61/63 (97%)	44 (72%)	13 (21%)	4 (7%)	1	19
All	All	11192/11458 (98%)	8966 (80%)	1699 (15%)	527 (5%)	3	29

All (527) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	AC	15	THR
4	AD	30	LYS
4	AD	137	SER
4	AD	138	TYR
4	AD	168	ARG
13	AM	63	THR

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Mol	Chain	Res	Type
13	AM	106	ASN
19	AS	28	LYS
20	AT	71	THR
25	BC	26	LYS
25	BC	33	LEU
25	BC	35	LYS
25	BC	237	GLU
25	BC	239	ARG
25	BC	260	ARG
26	BD	16	ARG
27	BE	73	ALA
28	BF	87	PRO
29	BG	92	ILE
29	BG	165	ALA
30	BH	91	SER
32	BJ	116	THR
32	BJ	149	PRO
32	BJ	157	ARG
34	BL	15	ARG
34	BL	36	LYS
34	BL	46	LYS
34	BL	59	LEU
34	BL	141	ALA
35	BM	8	LYS
35	BM	21	THR
37	BO	12	PHE
37	BO	59	LYS
37	BO	62	LYS
37	BO	90	GLY
37	BO	91	PRO
38	BP	58	ASN
38	BP	115	ARG
40	BR	53	GLU
40	BR	78	LYS
41	BS	110	LYS
43	BU	3	VAL
43	BU	7	VAL
44	BV	178	GLU
45	BW	47	PRO
46	BX	11	ARG
47	BY	47	ASN
50	B2	4	HIS

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Mol	Chain	Res	Type
50	B2	35	GLU
51	B3	28	ARG
3	CC	15	THR
4	CD	30	LYS
4	CD	137	SER
4	CD	138	TYR
4	CD	168	ARG
13	CM	63	THR
13	CM	106	ASN
19	CS	28	LYS
20	CT	71	THR
25	DC	26	LYS
25	DC	33	LEU
25	DC	35	LYS
25	DC	237	GLU
25	DC	239	ARG
25	DC	260	ARG
26	DD	16	ARG
27	DE	73	ALA
28	DF	87	PRO
29	DG	92	ILE
29	DG	165	ALA
30	DH	91	SER
32	DJ	116	THR
32	DJ	149	PRO
32	DJ	157	ARG
34	DL	15	ARG
34	DL	36	LYS
34	DL	46	LYS
34	DL	59	LEU
34	DL	141	ALA
35	DM	8	LYS
35	DM	21	THR
37	DO	12	PHE
37	DO	59	LYS
37	DO	62	LYS
37	DO	90	GLY
37	DO	91	PRO
38	DP	58	ASN
38	DP	115	ARG
40	DR	53	GLU
40	DR	78	LYS

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
41	DS	110	LYS
43	DU	3	VAL
43	DU	7	VAL
44	DV	178	GLU
45	DW	47	PRO
46	DX	11	ARG
47	DY	47	ASN
50	D2	35	GLU
51	D3	28	ARG
2	AB	18	GLY
2	AB	150	SER
4	AD	171	GLY
10	AJ	92	THR
12	AL	45	LYS
12	AL	63	TYR
13	AM	4	ILE
14	AN	26	ARG
19	AS	11	VAL
19	AS	29	ARG
20	AT	9	ASN
25	BC	34	VAL
25	BC	69	ARG
25	BC	70	TRP
25	BC	106	ILE
25	BC	197	GLY
26	BD	86	PRO
27	BE	84	VAL
28	BF	14	GLU
28	BF	24	GLY
28	BF	86	MET
30	BH	10	GLU
30	BH	90	GLY
30	BH	143	SER
32	BJ	106	LYS
32	BJ	153	HIS
34	BL	11	GLY
34	BL	34	GLY
34	BL	149	GLU
35	BM	7	MET
35	BM	10	ARG
35	BM	18	LYS
35	BM	62	GLY

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Mol	Chain	Res	Type
36	BN	3	HIS
36	BN	57	ARG
37	BO	35	ILE
37	BO	44	LYS
37	BO	57	LYS
38	BP	2	ASN
43	BU	17	SER
43	BU	42	VAL
43	BU	80	GLY
43	BU	98	VAL
44	BV	114	GLY
44	BV	120	ILE
44	BV	177	PRO
44	BV	179	ASP
45	BW	13	GLY
45	BW	73	GLY
46	BX	85	LEU
53	B5	3	LYS
2	CB	18	GLY
2	CB	150	SER
4	CD	171	GLY
10	CJ	92	THR
12	CL	45	LYS
12	CL	63	TYR
13	CM	4	ILE
14	CN	26	ARG
19	CS	11	VAL
19	CS	29	ARG
20	CT	9	ASN
25	DC	34	VAL
25	DC	69	ARG
25	DC	70	TRP
25	DC	106	ILE
25	DC	125	ILE
25	DC	197	GLY
26	DD	86	PRO
27	DE	84	VAL
28	DF	14	GLU
28	DF	24	GLY
28	DF	86	MET
30	DH	10	GLU
30	DH	90	GLY

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Mol	Chain	Res	Type
32	DJ	106	LYS
32	DJ	153	HIS
34	DL	11	GLY
34	DL	34	GLY
34	DL	149	GLU
35	DM	7	MET
35	DM	10	ARG
35	DM	18	LYS
35	DM	62	GLY
36	DN	3	HIS
36	DN	57	ARG
37	DO	35	ILE
37	DO	44	LYS
37	DO	57	LYS
38	DP	2	ASN
43	DU	17	SER
43	DU	42	VAL
43	DU	80	GLY
43	DU	98	VAL
44	DV	120	ILE
44	DV	177	PRO
44	DV	179	ASP
45	DW	13	GLY
45	DW	73	GLY
46	DX	85	LEU
50	D2	4	HIS
53	D5	3	LYS
2	AB	19	HIS
3	AC	45	LYS
3	AC	47	LEU
3	AC	105	GLU
3	AC	189	ALA
4	AD	32	ALA
4	AD	40	PRO
6	AF	49	ALA
7	AG	7	ALA
10	AJ	32	ALA
11	AK	49	GLY
12	AL	50	ALA
13	AM	101	GLN
19	AS	31	ILE
20	AT	98	PRO

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Mol	Chain	Res	Type
21	AU	9	ARG
25	BC	125	ILE
25	BC	238	GLY
26	BD	18	ASP
27	BE	68	LYS
27	BE	127	GLU
27	BE	166	ALA
29	BG	21	PRO
32	BJ	148	GLY
33	BK	4	PRO
33	BK	26	LYS
33	BK	97	ARG
34	BL	18	ARG
34	BL	42	SER
34	BL	49	ARG
34	BL	148	LEU
35	BM	133	ARG
36	BN	58	GLY
37	BO	95	HIS
38	BP	36	GLU
40	BR	17	GLY
40	BR	29	PRO
40	BR	80	GLN
43	BU	88	LYS
43	BU	96	ILE
44	BV	168	GLU
46	BX	9	GLY
46	BX	31	GLY
46	BX	32	LYS
47	BY	17	SER
49	B1	44	CYS
49	B1	62	CYS
51	B3	31	PRO
51	B3	46	HIS
53	B5	34	TRP
53	B5	35	GLN
2	CB	19	HIS
3	CC	47	LEU
3	CC	105	GLU
3	CC	189	ALA
4	CD	32	ALA
4	CD	40	PRO

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Mol	Chain	Res	Type
6	CF	49	ALA
7	CG	7	ALA
10	CJ	32	ALA
10	CJ	57	LYS
11	CK	49	GLY
13	CM	101	GLN
16	CP	48	TRP
19	CS	31	ILE
20	CT	98	PRO
21	CU	9	ARG
26	DD	18	ASP
27	DE	68	LYS
27	DE	166	ALA
29	DG	21	PRO
30	DH	143	SER
32	DJ	148	GLY
33	DK	4	PRO
33	DK	26	LYS
33	DK	97	ARG
34	DL	18	ARG
34	DL	42	SER
34	DL	49	ARG
34	DL	148	LEU
36	DN	58	GLY
37	DO	85	VAL
37	DO	95	HIS
38	DP	36	GLU
38	DP	57	PHE
40	DR	17	GLY
40	DR	29	PRO
40	DR	80	GLN
43	DU	88	LYS
43	DU	96	ILE
44	DV	114	GLY
44	DV	168	GLU
46	DX	9	GLY
46	DX	31	GLY
46	DX	32	LYS
47	DY	17	SER
49	D1	44	CYS
49	D1	62	CYS
51	D3	31	PRO

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Mol	Chain	Res	Type
51	D3	46	HIS
53	D5	34	TRP
53	D5	35	GLN
2	AB	130	ARG
3	AC	60	ALA
3	AC	129	ALA
5	AE	38	GLN
5	AE	72	GLN
9	AI	24	GLY
9	AI	58	ARG
10	AJ	57	LYS
11	AK	90	GLY
12	AL	64	GLU
14	AN	18	VAL
17	AQ	11	VAL
18	AR	36	ASN
19	AS	27	GLU
25	BC	32	SER
25	BC	191	ALA
25	BC	256	GLY
26	BD	43	GLY
28	BF	35	GLU
28	BF	124	SER
30	BH	132	PRO
32	BJ	70	ALA
33	BK	29	ASN
34	BL	65	ARG
35	BM	81	VAL
35	BM	134	ARG
35	BM	140	ALA
36	BN	8	ARG
37	BO	83	LYS
37	BO	85	VAL
37	BO	101	LEU
38	BP	57	PHE
39	BQ	91	ASP
40	BR	2	PHE
41	BS	11	ARG
42	BT	4	ALA
43	BU	39	VAL
43	BU	90	LEU
44	BV	80	ARG

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Mol	Chain	Res	Type
44	BV	142	SER
47	BY	15	LYS
48	BZ	29	ARG
49	B1	54	LYS
50	B2	45	VAL
51	B3	32	ASN
51	B3	51	GLU
53	B5	20	GLY
2	CB	130	ARG
3	CC	45	LYS
3	CC	60	ALA
3	CC	129	ALA
5	CE	38	GLN
5	CE	72	GLN
9	CI	24	GLY
9	CI	58	ARG
11	CK	90	GLY
12	CL	50	ALA
12	CL	64	GLU
14	CN	18	VAL
17	CQ	11	VAL
18	CR	36	ASN
19	CS	27	GLU
25	DC	32	SER
25	DC	238	GLY
25	DC	256	GLY
26	DD	43	GLY
27	DE	127	GLU
28	DF	35	GLU
28	DF	124	SER
30	DH	132	PRO
32	DJ	70	ALA
33	DK	29	ASN
34	DL	65	ARG
35	DM	81	VAL
35	DM	133	ARG
35	DM	134	ARG
35	DM	140	ALA
36	DN	8	ARG
37	DO	83	LYS
37	DO	101	LEU
39	DQ	91	ASP

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Mol	Chain	Res	Type
40	DR	2	PHE
41	DS	11	ARG
42	DT	4	ALA
43	DU	39	VAL
43	DU	90	LEU
44	DV	80	ARG
44	DV	142	SER
47	DY	15	LYS
48	DZ	29	ARG
49	D1	54	LYS
50	D2	45	VAL
51	D3	32	ASN
51	D3	51	GLU
53	D5	20	GLY
2	AB	15	VAL
2	AB	234	PRO
3	AC	127	ARG
3	AC	159	GLY
5	AE	49	PRO
5	AE	64	ARG
9	AI	100	GLY
10	AJ	54	PHE
11	AK	118	GLY
11	AK	125	PHE
12	AL	82	VAL
15	AO	23	GLY
16	AP	48	TRP
18	AR	78	LEU
25	BC	198	ASN
28	BF	12	TYR
35	BM	25	ASP
37	BO	61	ASN
38	BP	55	ASN
40	BR	48	GLY
46	BX	53	VAL
46	BX	84	GLY
46	BX	87	PRO
53	B5	61	LEU
2	CB	15	VAL
2	CB	234	PRO
3	CC	81	GLY
3	CC	127	ARG

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Mol	Chain	Res	Type
3	CC	159	GLY
5	CE	49	PRO
9	CI	100	GLY
10	CJ	54	PHE
11	CK	118	GLY
12	CL	82	VAL
15	CO	23	GLY
18	CR	78	LEU
25	DC	191	ALA
25	DC	198	ASN
28	DF	12	TYR
35	DM	25	ASP
37	DO	61	ASN
38	DP	55	ASN
40	DR	48	GLY
46	DX	53	VAL
46	DX	84	GLY
46	DX	87	PRO
3	AC	81	GLY
3	AC	96	GLY
3	AC	145	GLY
10	AJ	75	ILE
14	AN	14	PRO
26	BD	29	GLY
28	BF	142	PRO
29	BG	39	PRO
34	BL	43	GLY
35	BM	27	VAL
36	BN	13	HIS
43	BU	11	ASP
46	BX	58	ILE
3	CC	96	GLY
3	CC	145	GLY
4	CD	86	LYS
5	CE	64	ARG
10	CJ	75	ILE
11	CK	125	PHE
14	CN	14	PRO
16	CP	25	ARG
18	CR	20	ALA
26	DD	29	GLY
28	DF	136	ARG

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Mol	Chain	Res	Type
28	DF	142	PRO
29	DG	39	PRO
34	DL	43	GLY
35	DM	27	VAL
43	DU	11	ASP
3	AC	14	ILE
7	AG	130	GLY
15	AO	86	GLY
25	BC	236	GLY
34	BL	10	PRO
35	BM	96	VAL
3	CC	14	ILE
7	CG	130	GLY
15	CO	86	GLY
25	DC	236	GLY
33	DK	27	GLY
34	DL	10	PRO
46	DX	58	ILE
7	AG	82	GLY
12	AL	28	GLY
16	AP	63	GLY
27	BE	131	GLY
33	BK	27	GLY
40	BR	61	VAL
7	CG	82	GLY
12	CL	28	GLY
16	CP	63	GLY
27	DE	131	GLY
35	DM	96	VAL
40	DR	61	VAL
26	BD	4	ILE
35	BM	30	GLY
40	BR	16	PRO
44	BV	39	VAL
26	DD	4	ILE
33	DK	101	PRO
39	DQ	90	VAL
40	DR	16	PRO
44	DV	101	PRO
4	AD	146	ILE
10	AJ	77	PRO
26	BD	189	PRO

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Mol	Chain	Res	Type
27	BE	82	ILE
30	BH	144	VAL
33	BK	101	PRO
33	BK	119	PRO
39	BQ	90	VAL
44	BV	101	PRO
4	CD	146	ILE
32	DJ	158	PRO
33	DK	119	PRO
44	DV	39	VAL
50	D2	46	CYS
26	BD	173	VAL
32	BJ	158	PRO
10	CJ	77	PRO
27	DE	82	ILE
30	DH	144	VAL
43	DU	41	GLY

### 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	AB	202/202 (100%)	192 (95%)	10 (5%)	30	69
2	CB	202/202 (100%)	192 (95%)	10 (5%)	30	69
3	AC	160/160 (100%)	154 (96%)	6 (4%)	40	76
3	CC	160/160 (100%)	154 (96%)	6 (4%)	40	76
4	AD	180/180 (100%)	168 (93%)	12 (7%)	20	61
4	CD	180/180 (100%)	167 (93%)	13 (7%)	18	57
5	AE	116/116 (100%)	108 (93%)	8 (7%)	19	59
5	CE	116/116 (100%)	108 (93%)	8 (7%)	19	59
6	AF	90/90 (100%)	85 (94%)	5 (6%)	26	66
6	CF	90/90 (100%)	85 (94%)	5 (6%)	26	66
7	AG	126/126 (100%)	126 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
7	CG	126/126 (100%)	126 (100%)	0	100	100
8	AH	119/119 (100%)	114 (96%)	5 (4%)	36	74
8	CH	119/119 (100%)	114 (96%)	5 (4%)	36	74
9	AI	98/98 (100%)	92 (94%)	6 (6%)	23	64
9	CI	98/98 (100%)	92 (94%)	6 (6%)	23	64
10	AJ	88/88 (100%)	79 (90%)	9 (10%)	9	40
10	CJ	88/88 (100%)	79 (90%)	9 (10%)	9	40
11	AK	90/90 (100%)	86 (96%)	4 (4%)	35	73
11	CK	90/90 (100%)	86 (96%)	4 (4%)	35	73
12	AL	104/104 (100%)	94 (90%)	10 (10%)	10	43
12	CL	104/104 (100%)	94 (90%)	10 (10%)	10	43
13	AM	94/94 (100%)	88 (94%)	6 (6%)	22	62
13	CM	94/94 (100%)	88 (94%)	6 (6%)	22	62
14	AN	49/49 (100%)	47 (96%)	2 (4%)	37	74
14	CN	49/49 (100%)	48 (98%)	1 (2%)	63	87
15	AO	79/79 (100%)	75 (95%)	4 (5%)	29	69
15	CO	79/79 (100%)	75 (95%)	4 (5%)	29	69
16	AP	72/72 (100%)	67 (93%)	5 (7%)	19	59
16	CP	72/72 (100%)	68 (94%)	4 (6%)	26	66
17	AQ	94/94 (100%)	92 (98%)	2 (2%)	61	86
17	CQ	94/94 (100%)	92 (98%)	2 (2%)	61	86
18	AR	61/61 (100%)	59 (97%)	2 (3%)	45	79
18	CR	61/61 (100%)	59 (97%)	2 (3%)	45	79
19	AS	69/69 (100%)	60 (87%)	9 (13%)	5	27
19	CS	69/69 (100%)	60 (87%)	9 (13%)	5	27
20	AT	76/76 (100%)	71 (93%)	5 (7%)	21	61
20	CT	76/76 (100%)	71 (93%)	5 (7%)	21	61
21	AU	19/19 (100%)	18 (95%)	1 (5%)	28	67
21	CU	19/19 (100%)	18 (95%)	1 (5%)	28	67
25	BC	213/213 (100%)	192 (90%)	21 (10%)	10	41
25	DC	213/213 (100%)	192 (90%)	21 (10%)	10	41

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
26	BD	165/165 (100%)	149 (90%)	16 (10%)	10	42
26	DD	165/165 (100%)	149 (90%)	16 (10%)	10	42
27	BE	161/161 (100%)	148 (92%)	13 (8%)	15	52
27	DE	161/161 (100%)	147 (91%)	14 (9%)	13	48
28	BF	155/155 (100%)	144 (93%)	11 (7%)	18	58
28	DF	155/155 (100%)	144 (93%)	11 (7%)	18	58
29	BG	132/132 (100%)	123 (93%)	9 (7%)	20	60
29	DG	132/132 (100%)	123 (93%)	9 (7%)	20	60
30	BH	122/122 (100%)	115 (94%)	7 (6%)	25	66
30	DH	122/122 (100%)	115 (94%)	7 (6%)	25	66
31	BI	27/53 (51%)	26 (96%)	1 (4%)	41	76
31	DI	27/53 (51%)	26 (96%)	1 (4%)	41	76
32	BJ	116/116 (100%)	103 (89%)	13 (11%)	7	35
32	DJ	116/116 (100%)	103 (89%)	13 (11%)	7	35
33	BK	100/100 (100%)	92 (92%)	8 (8%)	15	52
33	DK	100/100 (100%)	92 (92%)	8 (8%)	15	52
34	BL	112/112 (100%)	87 (78%)	25 (22%)	1	6
34	DL	112/112 (100%)	87 (78%)	25 (22%)	1	6
35	BM	106/106 (100%)	98 (92%)	8 (8%)	17	55
35	DM	106/106 (100%)	98 (92%)	8 (8%)	17	55
36	BN	100/100 (100%)	94 (94%)	6 (6%)	24	64
36	DN	100/100 (100%)	94 (94%)	6 (6%)	24	64
37	BO	77/77 (100%)	68 (88%)	9 (12%)	7	32
37	DO	77/77 (100%)	68 (88%)	9 (12%)	7	32
38	BP	121/121 (100%)	109 (90%)	12 (10%)	10	41
38	DP	121/121 (100%)	110 (91%)	11 (9%)	12	46
39	BQ	92/92 (100%)	88 (96%)	4 (4%)	35	74
39	DQ	92/92 (100%)	88 (96%)	4 (4%)	35	74
40	BR	82/82 (100%)	77 (94%)	5 (6%)	23	64
40	DR	82/82 (100%)	77 (94%)	5 (6%)	23	64
41	BS	91/91 (100%)	85 (93%)	6 (7%)	21	61

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
41	DS	91/91 (100%)	85 (93%)	6 (7%)	21	61
42	BT	74/74 (100%)	67 (90%)	7 (10%)	11	43
42	DT	74/74 (100%)	67 (90%)	7 (10%)	11	43
43	BU	84/84 (100%)	78 (93%)	6 (7%)	18	58
43	DU	84/84 (100%)	78 (93%)	6 (7%)	18	58
44	BV	163/163 (100%)	159 (98%)	4 (2%)	55	84
44	DV	163/163 (100%)	159 (98%)	4 (2%)	55	84
45	BW	61/61 (100%)	55 (90%)	6 (10%)	10	42
45	DW	61/61 (100%)	55 (90%)	6 (10%)	10	42
46	BX	73/73 (100%)	61 (84%)	12 (16%)	3	17
46	DX	73/73 (100%)	61 (84%)	12 (16%)	3	17
47	BY	58/58 (100%)	51 (88%)	7 (12%)	6	30
47	DY	58/58 (100%)	51 (88%)	7 (12%)	6	30
48	BZ	51/51 (100%)	49 (96%)	2 (4%)	39	76
48	DZ	51/51 (100%)	49 (96%)	2 (4%)	39	76
49	B1	27/27 (100%)	24 (89%)	3 (11%)	8	35
49	D1	27/27 (100%)	24 (89%)	3 (11%)	8	35
50	B2	45/45 (100%)	43 (96%)	2 (4%)	35	73
50	D2	45/45 (100%)	43 (96%)	2 (4%)	35	73
51	B3	43/43 (100%)	39 (91%)	4 (9%)	11	45
51	D3	43/43 (100%)	39 (91%)	4 (9%)	11	45
52	B4	41/41 (100%)	34 (83%)	7 (17%)	2	15
52	D4	41/41 (100%)	34 (83%)	7 (17%)	2	15
53	B5	53/53 (100%)	51 (96%)	2 (4%)	40	76
53	D5	53/53 (100%)	51 (96%)	2 (4%)	40	76
All	All	9462/9514 (100%)	8769 (93%)	693 (7%)	17	57

All (693) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	AB	25	ASN
2	AB	27	LYS
2	AB	71	VAL

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Mol	Chain	Res	Type
2	AB	75	LYS
2	AB	116	GLU
2	AB	117	GLU
2	AB	153	ARG
2	AB	154	LEU
2	AB	178	ARG
2	AB	221	LEU
3	AC	3	ASN
3	AC	5	ILE
3	AC	27	LYS
3	AC	79	ARG
3	AC	91	LEU
3	AC	196	LEU
4	AD	3	ARG
4	AD	21	LEU
4	AD	49	ARG
4	AD	72	GLU
4	AD	73	ARG
4	AD	119	GLN
4	AD	122	ARG
4	AD	131	ARG
4	AD	135	LEU
4	AD	150	GLU
4	AD	158	ILE
4	AD	166	LYS
5	AE	8	GLU
5	AE	12	LEU
5	AE	16	THR
5	AE	47	LYS
5	AE	76	ILE
5	AE	79	GLU
5	AE	137	GLU
5	AE	144	THR
6	AF	48	LEU
6	AF	59	TYR
6	AF	78	GLU
6	AF	98	LEU
6	AF	100	ASN
8	AH	1	MET
8	AH	25	ASP
8	AH	30	ARG
8	AH	102	ARG

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Mol	Chain	Res	Type
8	AH	136	GLU
9	AI	10	ARG
9	AI	19	LEU
9	AI	95	LYS
9	AI	99	LEU
9	AI	104	ARG
9	AI	121	ARG
10	AJ	16	LEU
10	AJ	22	LYS
10	AJ	49	VAL
10	AJ	55	LYS
10	AJ	62	HIS
10	AJ	74	ILE
10	AJ	80	LYS
10	AJ	92	THR
10	AJ	96	ILE
11	AK	33	THR
11	AK	92	GLU
11	AK	117	ASN
11	AK	119	CYS
12	AL	6	ILE
12	AL	19	LYS
12	AL	26	LEU
12	AL	37	THR
12	AL	40	ARG
12	AL	41	THR
12	AL	64	GLU
12	AL	65	VAL
12	AL	76	LEU
12	AL	81	VAL
13	AM	58	GLU
13	AM	64	TRP
13	AM	87	TYR
13	AM	93	ARG
13	AM	106	ASN
13	AM	115	LYS
14	AN	16	PHE
14	AN	40	CYS
15	AO	5	LYS
15	AO	17	ARG
15	AO	44	LYS
15	AO	82	ILE

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Mol	Chain	Res	Type
16	AP	2	VAL
16	AP	27	LYS
16	AP	67	THR
16	AP	82	GLN
16	AP	83	GLU
17	AQ	38	ARG
17	AQ	52	LYS
18	AR	42	ARG
18	AR	84	LYS
19	AS	5	LEU
19	AS	6	LYS
19	AS	7	LYS
19	AS	27	GLU
19	AS	29	ARG
19	AS	37	ARG
19	AS	44	MET
19	AS	53	ASN
19	AS	70	LYS
20	AT	22	ARG
20	AT	26	ASN
20	AT	60	GLU
20	AT	62	LEU
20	AT	93	GLU
21	AU	5	ASP
25	BC	5	LYS
25	BC	10	THR
25	BC	28	GLU
25	BC	33	LEU
25	BC	44	ASN
25	BC	50	THR
25	BC	61	LEU
25	BC	78	LYS
25	BC	94	LEU
25	BC	95	LEU
25	BC	106	ILE
25	BC	111	LEU
25	BC	150	LYS
25	BC	166	GLN
25	BC	192	THR
25	BC	237	GLU
25	BC	242	ARG
25	BC	244	ARG

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Mol	Chain	Res	Type
25	BC	259	THR
25	BC	261	LYS
25	BC	267	SER
26	BD	9	VAL
26	BD	48	GLN
26	BD	52	LEU
26	BD	54	GLN
26	BD	57	LYS
26	BD	92	THR
26	BD	118	LYS
26	BD	119	ARG
26	BD	132	HIS
26	BD	141	ILE
26	BD	144	ARG
26	BD	154	LYS
26	BD	160	TYR
26	BD	169	ASN
26	BD	184	VAL
26	BD	195	LEU
27	BE	8	GLN
27	BE	9	ILE
27	BE	54	ARG
27	BE	56	GLU
27	BE	66	PRO
27	BE	69	HIS
27	BE	72	ARG
27	BE	83	PHE
27	BE	95	ARG
27	BE	106	ARG
27	BE	164	ARG
27	BE	188	ARG
27	BE	195	ASP
28	BF	18	GLU
28	BF	33	ARG
28	BF	34	LEU
28	BF	47	LYS
28	BF	76	SER
28	BF	86	MET
28	BF	90	LEU
28	BF	98	ARG
28	BF	107	LEU
28	BF	115	ARG

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Mol	Chain	Res	Type
28	BF	155	MET
29	BG	13	LYS
29	BG	23	ARG
29	BG	43	VAL
29	BG	86	GLU
29	BG	101	ARG
29	BG	105	LEU
29	BG	123	PHE
29	BG	158	HIS
29	BG	162	ILE
30	BH	5	LEU
30	BH	6	LEU
30	BH	67	ARG
30	BH	73	GLU
30	BH	77	LEU
30	BH	92	VAL
30	BH	109	ILE
31	BI	3	ASN
32	BJ	57	LEU
32	BJ	58	ARG
32	BJ	64	ASP
32	BJ	71	MET
32	BJ	92	GLN
32	BJ	94	ILE
32	BJ	96	THR
32	BJ	105	LEU
32	BJ	110	LEU
32	BJ	117	HIS
32	BJ	120	ARG
32	BJ	135	LEU
32	BJ	161	LEU
33	BK	19	ILE
33	BK	25	LEU
33	BK	47	ILE
33	BK	77	ILE
33	BK	87	ILE
33	BK	99	PHE
33	BK	104	ARG
33	BK	122	LEU
34	BL	6	LEU
34	BL	13	ASN
34	BL	15	ARG

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Mol	Chain	Res	Type
34	BL	16	ARG
34	BL	19	VAL
34	BL	32	THR
34	BL	35	HIS
34	BL	38	GLN
34	BL	39	LYS
34	BL	49	ARG
34	BL	50	ARG
34	BL	52	GLU
34	BL	57	THR
34	BL	61	ARG
34	BL	62	LEU
34	BL	67	MET
34	BL	75	ILE
34	BL	81	GLN
34	BL	83	VAL
34	BL	84	ASN
34	BL	105	LEU
34	BL	106	LEU
34	BL	111	ARG
34	BL	147	LEU
34	BL	148	LEU
35	BM	6	ARG
35	BM	13	GLN
35	BM	14	ARG
35	BM	22	LYS
35	BM	45	GLN
35	BM	60	ARG
35	BM	89	ASN
35	BM	135	ASP
36	BN	5	LYS
36	BN	8	ARG
36	BN	9	LYS
36	BN	10	LEU
36	BN	79	LEU
36	BN	104	ARG
37	BO	18	ILE
37	BO	26	LEU
37	BO	30	ARG
37	BO	36	TYR
37	BO	42	ASP
37	BO	44	LYS

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
37	BO	61	ASN
37	BO	62	LYS
37	BO	93	LYS
38	BP	41	ARG
38	BP	58	ASN
38	BP	59	THR
38	BP	78	LEU
38	BP	86	ILE
38	BP	87	ASP
38	BP	89	VAL
38	BP	98	LYS
38	BP	99	LEU
38	BP	108	ARG
38	BP	112	ARG
38	BP	113	LYS
39	BQ	79	PHE
39	BQ	92	ARG
39	BQ	97	ASP
39	BQ	103	PRO
40	BR	12	TYR
40	BR	13	ARG
40	BR	18	LEU
40	BR	80	GLN
40	BR	99	ILE
41	BS	11	ARG
41	BS	23	LEU
41	BS	69	LEU
41	BS	70	TYR
41	BS	77	ASP
41	BS	107	LEU
42	BT	28	PHE
42	BT	57	LEU
42	BT	65	ARG
42	BT	68	ARG
42	BT	75	ASP
42	BT	80	ILE
42	BT	83	VAL
43	BU	4	LYS
43	BU	6	HIS
43	BU	8	LYS
43	BU	31	LEU
43	BU	76	CYS

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
43	BU	97	ARG
44	BV	25	PRO
44	BV	72	ARG
44	BV	76	LEU
44	BV	94	GLU
45	BW	14	ARG
45	BW	21	LEU
45	BW	25	ARG
45	BW	35	ASN
45	BW	64	ASP
45	BW	80	HIS
46	BX	11	ARG
46	BX	17	SER
46	BX	18	ILE
46	BX	20	ARG
46	BX	40	ARG
46	BX	45	ASN
46	BX	46	LEU
46	BX	73	LEU
46	BX	75	GLU
46	BX	82	LEU
46	BX	89	GLU
46	BX	95	LEU
47	BY	2	LYS
47	BY	21	LEU
47	BY	37	PHE
47	BY	53	LEU
47	BY	56	GLN
47	BY	59	ARG
47	BY	61	LEU
48	BZ	10	LYS
48	BZ	46	ASN
49	B1	46	ASN
49	B1	49	GLU
49	B1	60	GLU
50	B2	3	LYS
50	B2	23	HIS
51	B3	11	LEU
51	B3	29	ASN
51	B3	34	LEU
51	B3	42	TRP
52	B4	4	THR

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Mol	Chain	Res	Type
52	B4	8	ASN
52	B4	10	ARG
52	B4	15	THR
52	B4	19	ARG
52	B4	24	THR
52	B4	29	LYS
53	B5	4	MET
53	B5	30	ARG
2	CB	25	ASN
2	CB	27	LYS
2	CB	71	VAL
2	CB	75	LYS
2	CB	116	GLU
2	CB	117	GLU
2	CB	153	ARG
2	CB	154	LEU
2	CB	178	ARG
2	CB	221	LEU
3	CC	3	ASN
3	CC	5	ILE
3	CC	27	LYS
3	CC	79	ARG
3	CC	91	LEU
3	CC	196	LEU
4	CD	3	ARG
4	CD	21	LEU
4	CD	49	ARG
4	CD	72	GLU
4	CD	73	ARG
4	CD	98	GLU
4	CD	119	GLN
4	CD	122	ARG
4	CD	131	ARG
4	CD	135	LEU
4	CD	150	GLU
4	CD	158	ILE
4	CD	166	LYS
5	CE	8	GLU
5	CE	12	LEU
5	CE	16	THR
5	CE	47	LYS
5	CE	76	ILE

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Mol	Chain	Res	Type
5	CE	79	GLU
5	CE	137	GLU
5	CE	144	THR
6	CF	48	LEU
6	CF	59	TYR
6	CF	78	GLU
6	CF	98	LEU
6	CF	100	ASN
8	CH	1	MET
8	CH	25	ASP
8	CH	30	ARG
8	CH	102	ARG
8	CH	136	GLU
9	CI	10	ARG
9	CI	19	LEU
9	CI	95	LYS
9	CI	99	LEU
9	CI	104	ARG
9	CI	121	ARG
10	CJ	16	LEU
10	CJ	22	LYS
10	CJ	49	VAL
10	CJ	55	LYS
10	CJ	62	HIS
10	CJ	74	ILE
10	CJ	80	LYS
10	CJ	92	THR
10	CJ	96	ILE
11	CK	33	THR
11	CK	92	GLU
11	CK	117	ASN
11	CK	119	CYS
12	CL	6	ILE
12	CL	19	LYS
12	CL	26	LEU
12	CL	37	THR
12	CL	40	ARG
12	CL	41	THR
12	CL	64	GLU
12	CL	65	VAL
12	CL	76	LEU
12	CL	81	VAL

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Mol	Chain	Res	Type
13	CM	58	GLU
13	CM	64	TRP
13	CM	87	TYR
13	CM	93	ARG
13	CM	106	ASN
13	CM	115	LYS
14	CN	16	PHE
15	CO	5	LYS
15	CO	17	ARG
15	CO	44	LYS
15	CO	82	ILE
16	CP	2	VAL
16	CP	27	LYS
16	CP	82	GLN
16	CP	83	GLU
17	CQ	38	ARG
17	CQ	52	LYS
18	CR	42	ARG
18	CR	84	LYS
19	CS	5	LEU
19	CS	6	LYS
19	CS	7	LYS
19	CS	27	GLU
19	CS	29	ARG
19	CS	37	ARG
19	CS	44	MET
19	CS	53	ASN
19	CS	70	LYS
20	CT	22	ARG
20	CT	26	ASN
20	CT	60	GLU
20	CT	62	LEU
20	CT	93	GLU
21	CU	5	ASP
25	DC	5	LYS
25	DC	10	THR
25	DC	28	GLU
25	DC	33	LEU
25	DC	44	ASN
25	DC	50	THR
25	DC	61	LEU
25	DC	78	LYS

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Mol	Chain	Res	Type
25	DC	94	LEU
25	DC	95	LEU
25	DC	106	ILE
25	DC	111	LEU
25	DC	150	LYS
25	DC	166	GLN
25	DC	192	THR
25	DC	237	GLU
25	DC	242	ARG
25	DC	244	ARG
25	DC	259	THR
25	DC	261	LYS
25	DC	267	SER
26	DD	9	VAL
26	DD	48	GLN
26	DD	52	LEU
26	DD	54	GLN
26	DD	57	LYS
26	DD	92	THR
26	DD	118	LYS
26	DD	119	ARG
26	DD	132	HIS
26	DD	141	ILE
26	DD	144	ARG
26	DD	154	LYS
26	DD	160	TYR
26	DD	169	ASN
26	DD	184	VAL
26	DD	195	LEU
27	DE	8	GLN
27	DE	9	ILE
27	DE	54	ARG
27	DE	56	GLU
27	DE	66	PRO
27	DE	69	HIS
27	DE	72	ARG
27	DE	78	ILE
27	DE	83	PHE
27	DE	95	ARG
27	DE	106	ARG
27	DE	164	ARG
27	DE	188	ARG

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Mol	Chain	Res	Type
27	DE	195	ASP
28	DF	18	GLU
28	DF	33	ARG
28	DF	34	LEU
28	DF	47	LYS
28	DF	76	SER
28	DF	86	MET
28	DF	90	LEU
28	DF	98	ARG
28	DF	107	LEU
28	DF	115	ARG
28	DF	155	MET
29	DG	13	LYS
29	DG	23	ARG
29	DG	43	VAL
29	DG	86	GLU
29	DG	101	ARG
29	DG	105	LEU
29	DG	123	PHE
29	DG	158	HIS
29	DG	162	ILE
30	DH	5	LEU
30	DH	6	LEU
30	DH	67	ARG
30	DH	73	GLU
30	DH	77	LEU
30	DH	92	VAL
30	DH	109	ILE
31	DI	3	ASN
32	DJ	57	LEU
32	DJ	58	ARG
32	DJ	64	ASP
32	DJ	71	MET
32	DJ	92	GLN
32	DJ	94	ILE
32	DJ	96	THR
32	DJ	105	LEU
32	DJ	110	LEU
32	DJ	117	HIS
32	DJ	120	ARG
32	DJ	135	LEU
32	DJ	161	LEU

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Mol	Chain	Res	Type
33	DK	19	ILE
33	DK	25	LEU
33	DK	47	ILE
33	DK	77	ILE
33	DK	87	ILE
33	DK	99	PHE
33	DK	104	ARG
33	DK	122	LEU
34	DL	6	LEU
34	DL	13	ASN
34	DL	15	ARG
34	DL	16	ARG
34	DL	32	THR
34	DL	35	HIS
34	DL	38	GLN
34	DL	39	LYS
34	DL	40	SER
34	DL	49	ARG
34	DL	50	ARG
34	DL	52	GLU
34	DL	57	THR
34	DL	61	ARG
34	DL	62	LEU
34	DL	67	MET
34	DL	75	ILE
34	DL	81	GLN
34	DL	83	VAL
34	DL	84	ASN
34	DL	105	LEU
34	DL	106	LEU
34	DL	111	ARG
34	DL	147	LEU
34	DL	148	LEU
35	DM	6	ARG
35	DM	13	GLN
35	DM	14	ARG
35	DM	22	LYS
35	DM	45	GLN
35	DM	60	ARG
35	DM	89	ASN
35	DM	135	ASP
36	DN	5	LYS

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
36	DN	8	ARG
36	DN	9	LYS
36	DN	10	LEU
36	DN	79	LEU
36	DN	104	ARG
37	DO	18	ILE
37	DO	26	LEU
37	DO	30	ARG
37	DO	36	TYR
37	DO	42	ASP
37	DO	44	LYS
37	DO	61	ASN
37	DO	62	LYS
37	DO	93	LYS
38	DP	41	ARG
38	DP	58	ASN
38	DP	59	THR
38	DP	78	LEU
38	DP	86	ILE
38	DP	87	ASP
38	DP	89	VAL
38	DP	98	LYS
38	DP	108	ARG
38	DP	112	ARG
38	DP	113	LYS
39	DQ	79	PHE
39	DQ	92	ARG
39	DQ	97	ASP
39	DQ	103	PRO
40	DR	12	TYR
40	DR	13	ARG
40	DR	18	LEU
40	DR	80	GLN
40	DR	99	ILE
41	DS	11	ARG
41	DS	23	LEU
41	DS	69	LEU
41	DS	70	TYR
41	DS	77	ASP
41	DS	107	LEU
42	DT	28	PHE
42	DT	57	LEU

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Mol	Chain	Res	Type
42	DT	65	ARG
42	DT	68	ARG
42	DT	75	ASP
42	DT	80	ILE
42	DT	83	VAL
43	DU	4	LYS
43	DU	6	HIS
43	DU	8	LYS
43	DU	31	LEU
43	DU	76	CYS
43	DU	97	ARG
44	DV	25	PRO
44	DV	72	ARG
44	DV	76	LEU
44	DV	94	GLU
45	DW	14	ARG
45	DW	21	LEU
45	DW	25	ARG
45	DW	35	ASN
45	DW	64	ASP
45	DW	80	HIS
46	DX	11	ARG
46	DX	17	SER
46	DX	18	ILE
46	DX	20	ARG
46	DX	40	ARG
46	DX	45	ASN
46	DX	46	LEU
46	DX	73	LEU
46	DX	75	GLU
46	DX	82	LEU
46	DX	89	GLU
46	DX	95	LEU
47	DY	2	LYS
47	DY	21	LEU
47	DY	37	PHE
47	DY	53	LEU
47	DY	56	GLN
47	DY	59	ARG
47	DY	61	LEU
48	DZ	10	LYS
48	DZ	46	ASN

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Mol	Chain	Res	Type
49	D1	46	ASN
49	D1	49	GLU
49	D1	60	GLU
50	D2	3	LYS
50	D2	23	HIS
51	D3	11	LEU
51	D3	29	ASN
51	D3	34	LEU
51	D3	42	TRP
52	D4	4	THR
52	D4	8	ASN
52	D4	10	ARG
52	D4	15	THR
52	D4	19	ARG
52	D4	24	THR
52	D4	29	LYS
53	D5	4	MET
53	D5	30	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (257) such sidechains are listed below:

Mol	Chain	Res	Type
2	AB	16	HIS
2	AB	25	ASN
2	AB	37	ASN
2	AB	146	GLN
2	AB	212	GLN
3	AC	28	GLN
3	AC	69	HIS
3	AC	170	GLN
3	AC	176	HIS
4	AD	77	ASN
4	AD	119	GLN
4	AD	123	HIS
5	AE	20	GLN
5	AE	73	ASN
5	AE	78	HIS
6	AF	27	GLN
6	AF	32	ASN
6	AF	100	ASN
7	AG	13	GLN
7	AG	28	ASN

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Mol	Chain	Res	Type
7	AG	84	ASN
7	AG	106	GLN
8	AH	82	HIS
9	AI	23	ASN
9	AI	73	GLN
9	AI	117	HIS
9	AI	124	GLN
10	AJ	68	HIS
10	AJ	78	ASN
10	AJ	84	GLN
11	AK	38	ASN
11	AK	117	ASN
12	AL	7	ASN
12	AL	48	ASN
12	AL	74	HIS
13	AM	101	GLN
15	AO	37	ASN
15	AO	46	HIS
15	AO	53	HIS
16	AP	13	HIS
16	AP	82	GLN
17	AQ	16	GLN
19	AS	14	HIS
19	AS	47	HIS
19	AS	53	ASN
19	AS	57	HIS
20	AT	26	ASN
25	BC	58	HIS
25	BC	87	ASN
25	BC	96	HIS
25	BC	116	GLN
25	BC	126	GLN
25	BC	166	GLN
25	BC	186	HIS
25	BC	198	ASN
25	BC	227	ASN
25	BC	233	HIS
26	BD	60	ASN
26	BD	66	HIS
26	BD	129	HIS
26	BD	143	ASN
26	BD	169	ASN

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
26	BD	192	ASN
27	BE	67	GLN
27	BE	69	HIS
27	BE	75	HIS
27	BE	169	ASN
28	BF	27	ASN
28	BF	58	GLN
28	BF	66	GLN
28	BF	108	ASN
28	BF	121	ASN
29	BG	143	GLN
29	BG	147	ASN
30	BH	133	HIS
31	BI	3	ASN
31	BI	6	ASN
32	BJ	79	ASN
32	BJ	151	HIS
32	BJ	154	GLN
33	BK	89	ASN
34	BL	13	ASN
34	BL	27	HIS
34	BL	35	HIS
34	BL	38	GLN
34	BL	81	GLN
35	BM	13	GLN
35	BM	45	GLN
35	BM	141	GLN
36	BN	13	HIS
36	BN	16	HIS
36	BN	61	HIS
36	BN	71	GLN
36	BN	91	GLN
37	BO	61	ASN
37	BO	95	HIS
38	BP	43	GLN
38	BP	58	ASN
38	BP	79	HIS
38	BP	84	GLN
38	BP	90	GLN
39	BQ	49	HIS
39	BQ	72	HIS
41	BS	34	ASN

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Mol	Chain	Res	Type
41	BS	57	ASN
41	BS	61	ASN
41	BS	102	HIS
42	BT	31	HIS
42	BT	41	ASN
42	BT	55	ASN
42	BT	87	GLN
43	BU	6	HIS
44	BV	118	GLN
45	BW	35	ASN
45	BW	50	ASN
45	BW	70	GLN
46	BX	19	GLN
46	BX	45	ASN
46	BX	56	GLN
46	BX	66	HIS
47	BY	47	ASN
48	BZ	19	GLN
48	BZ	46	ASN
48	BZ	52	HIS
49	B1	46	ASN
50	B2	43	HIS
51	B3	29	ASN
52	B4	8	ASN
52	B4	36	GLN
53	B5	31	HIS
53	B5	33	ASN
2	CB	16	HIS
2	CB	37	ASN
2	CB	146	GLN
2	CB	212	GLN
3	CC	28	GLN
3	CC	69	HIS
3	CC	170	GLN
4	CD	77	ASN
4	CD	119	GLN
4	CD	123	HIS
5	CE	20	GLN
5	CE	73	ASN
5	CE	78	HIS
6	CF	27	GLN
6	CF	32	ASN

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Mol	Chain	Res	Type
6	CF	100	ASN
7	CG	13	GLN
7	CG	84	ASN
7	CG	106	GLN
8	CH	78	GLN
8	CH	82	HIS
9	CI	23	ASN
9	CI	117	HIS
9	CI	124	GLN
10	CJ	68	HIS
10	CJ	78	ASN
10	CJ	84	GLN
11	CK	38	ASN
11	CK	117	ASN
12	CL	7	ASN
12	CL	74	HIS
13	CM	101	GLN
15	CO	37	ASN
15	CO	46	HIS
15	CO	53	HIS
16	CP	82	GLN
17	CQ	16	GLN
19	CS	14	HIS
19	CS	47	HIS
19	CS	53	ASN
19	CS	57	HIS
20	CT	26	ASN
25	DC	58	HIS
25	DC	87	ASN
25	DC	96	HIS
25	DC	116	GLN
25	DC	126	GLN
25	DC	166	GLN
25	DC	186	HIS
25	DC	198	ASN
25	DC	227	ASN
25	DC	233	HIS
26	DD	60	ASN
26	DD	66	HIS
26	DD	129	HIS
26	DD	143	ASN
26	DD	169	ASN

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Mol	Chain	Res	Type
26	DD	192	ASN
27	DE	67	GLN
27	DE	69	HIS
27	DE	75	HIS
27	DE	169	ASN
28	DF	27	ASN
28	DF	58	GLN
28	DF	66	GLN
28	DF	108	ASN
28	DF	121	ASN
29	DG	143	GLN
29	DG	147	ASN
30	DH	133	HIS
31	DI	3	ASN
31	DI	6	ASN
32	DJ	79	ASN
32	DJ	151	HIS
32	DJ	154	GLN
33	DK	89	ASN
34	DL	13	ASN
34	DL	27	HIS
34	DL	35	HIS
34	DL	38	GLN
34	DL	81	GLN
35	DM	13	GLN
35	DM	45	GLN
35	DM	141	GLN
36	DN	13	HIS
36	DN	16	HIS
36	DN	61	HIS
36	DN	71	GLN
36	DN	91	GLN
37	DO	61	ASN
37	DO	95	HIS
38	DP	43	GLN
38	DP	58	ASN
38	DP	79	HIS
38	DP	84	GLN
38	DP	90	GLN
39	DQ	49	HIS
39	DQ	72	HIS
41	DS	34	ASN

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Mol	Chain	Res	Type
41	DS	57	ASN
41	DS	61	ASN
41	DS	102	HIS
42	DT	31	HIS
42	DT	41	ASN
42	DT	55	ASN
42	DT	87	GLN
43	DU	6	HIS
44	DV	118	GLN
45	DW	35	ASN
45	DW	50	ASN
45	DW	70	GLN
46	DX	19	GLN
46	DX	45	ASN
46	DX	56	GLN
46	DX	66	HIS
47	DY	47	ASN
48	DZ	19	GLN
48	DZ	46	ASN
48	DZ	52	HIS
49	D1	46	ASN
50	D2	43	HIS
51	D3	29	ASN
52	D4	8	ASN
52	D4	36	GLN
53	D5	31	HIS
53	D5	33	ASN

### 5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	AA	1505/1506 (99%)	204 (13%)	35 (2%)
1	CA	1505/1506 (99%)	205 (13%)	35 (2%)
22	AV	34/35 (97%)	2 (5%)	2 (5%)
22	CV	34/35 (97%)	2 (5%)	2 (5%)
23	BA	2757/2879 (95%)	410 (14%)	69 (2%)
23	DA	2757/2879 (95%)	407 (14%)	70 (2%)
24	BB	118/119 (99%)	16 (13%)	1 (0%)
24	DB	118/119 (99%)	16 (13%)	1 (0%)
All	All	8828/9078 (97%)	1262 (14%)	215 (2%)

All (1262) RNA backbone outliers are listed below:



Mol	Chain	Res	Type
1	AA	9	G
1	AA	22	G
1	AA	32	A
1	AA	39	G
1	AA	41	G
1	AA	47	C
1	AA	48	C
1	AA	51	A
1	AA	61	G
1	AA	88	C
1	AA	99	C
1	AA	101	A
1	AA	108	G
1	AA	116	A
1	AA	120	A
1	AA	121	C
1	AA	131	C
1	AA	169	C
1	AA	182	U
1	AA	195	A
1	AA	197	A
1	AA	200	G
1	AA	209	U
1	AA	210	U
1	AA	243	A
1	AA	244	U
1	AA	247	G
1	AA	251	G
1	AA	266	G
1	AA	267	C
1	AA	289	G
1	AA	321	A
1	AA	328	C
1	AA	329	A
1	AA	332	G
1	AA	344	A
1	AA	345	C
1	AA	346	G
1	AA	352	C
1	AA	353	A
1	AA	354	G
1	AA	358	U
1	AA	367	U

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
1	AA	372	C
1	AA	373	A
1	AA	382	A
1	AA	387	U
1	AA	397	A
1	AA	398	C
1	AA	412	A
1	AA	413	G
1	AA	414	A
1	AA	422	C
1	AA	423	G
1	AA	429	U
1	AA	430	A
1	AA	439	A
1	AA	465	A
1	AA	467	G
1	AA	485	G
1	AA	496	A
1	AA	497	U
1	AA	511	C
1	AA	512	U
1	AA	518	C
1	AA	527	G
1	AA	531	U
1	AA	532	A
1	AA	533	A
1	AA	547	A
1	AA	559	A
1	AA	561	U
1	AA	562	C
1	AA	572	A
1	AA	573	A
1	AA	576	G
1	AA	596	C
1	AA	653	A
1	AA	687	A
1	AA	688	G
1	AA	703	G
1	AA	731	G
1	AA	755	G
1	AA	777	A
1	AA	793	U

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Mol	Chain	Res	Type
1	AA	794	A
1	AA	796	C
1	AA	816	A
1	AA	817	C
1	AA	818	G
1	AA	819	A
1	AA	828	A
1	AA	829	G
1	AA	833	U
1	AA	841	U
1	AA	842	C
1	AA	843	U
1	AA	855	G
1	AA	859	A
1	AA	873	A
1	AA	902	G
1	AA	914	A
1	AA	926	G
1	AA	927	G
1	AA	934	C
1	AA	935	A
1	AA	960	U
1	AA	961	U
1	AA	969	A
1	AA	971	G
1	AA	974	A
1	AA	976	G
1	AA	977	A
1	AA	980	C
1	AA	981	U
1	AA	992	U
1	AA	993	G
1	AA	1004	A
1	AA	1006	C
1	AA	1045	C
1	AA	1054	C
1	AA	1065	U
1	AA	1066	C
1	AA	1068	G
1	AA	1081	G
1	AA	1094	G
1	AA	1095	U

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Mol	Chain	Res	Type
1	AA	1101	A
1	AA	1113	C
1	AA	1117	G
1	AA	1118	C
1	AA	1125	U
1	AA	1126	U
1	AA	1129	C
1	AA	1130	A
1	AA	1131	G
1	AA	1136	U
1	AA	1137	C
1	AA	1139	G
1	AA	1146	A
1	AA	1151	A
1	AA	1152	A
1	AA	1159	U
1	AA	1160	G
1	AA	1181	G
1	AA	1182	G
1	AA	1193	G
1	AA	1196	U
1	AA	1197	G
1	AA	1202	G
1	AA	1212	U
1	AA	1213	A
1	AA	1214	C
1	AA	1224	G
1	AA	1225	A
1	AA	1238	A
1	AA	1240	U
1	AA	1241	G
1	AA	1256	A
1	AA	1257	U
1	AA	1280	A
1	AA	1281	U
1	AA	1282	C
1	AA	1285	A
1	AA	1286	A
1	AA	1287	A
1	AA	1300	G
1	AA	1301	U
1	AA	1302	U

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Mol	Chain	Res	Type
1	AA	1305	G
1	AA	1317	C
1	AA	1320	C
1	AA	1322	C
1	AA	1331	G
1	AA	1335	C
1	AA	1346	A
1	AA	1347	G
1	AA	136(B)	C
1	AA	1363	A
1	AA	1364	U
1	AA	1400	C
1	AA	1419	G
1	AA	1442	G
1	AA	1443	G
1	AA	1446	A
1	AA	1447	G
1	AA	1451	A
1	AA	1452	C
1	AA	1454	G
1	AA	1493	A
1	AA	1494	G
1	AA	1497	G
1	AA	1499	A
1	AA	1502	A
1	AA	1503	A
1	AA	1504	G
1	AA	1505	G
1	AA	1506	U
1	AA	1517	G
1	AA	1519	A
1	AA	1520	G
1	AA	1528	U
1	AA	1529	G
1	AA	1530	G
22	AV	6172	U
22	AV	6189	A
23	BA	10	G
23	BA	35	G
23	BA	46	C
23	BA	49	A
23	BA	64	A

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Mol	Chain	Res	Type
23	BA	71	A
23	BA	72	U
23	BA	74	A
23	BA	75	G
23	BA	84	A
23	BA	88	G
23	BA	102	G
23	BA	118	A
23	BA	119	A
23	BA	120	U
23	BA	125	G
23	BA	131	G
23	BA	138	G
23	BA	140	A
23	BA	181	A
23	BA	196	A
23	BA	197	A
23	BA	199	A
23	BA	205	G
23	BA	215	G
23	BA	216	A
23	BA	221	A
23	BA	222	A
23	BA	228	A
23	BA	229	A
23	BA	230	U
23	BA	245	G
23	BA	248	G
23	BA	252	G
23	BA	269	U
23	BA	270(K)	G
23	BA	270(M)	U
23	BA	270(N)	U
23	BA	270(O)	G
23	BA	270(R)	C
23	BA	271(D)	U
23	BA	271	G
23	BA	274	G
23	BA	275	G
23	BA	276	A
23	BA	277	C
23	BA	278	A

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
23	BA	279	C
23	BA	283	A
23	BA	302	C
23	BA	311	A
23	BA	323	G
23	BA	324	A
23	BA	329	G
23	BA	330	A
23	BA	332	A
23	BA	333	G
23	BA	343	C
23	BA	352	G
23	BA	353	G
23	BA	386	G
23	BA	396	G
23	BA	405	U
23	BA	411	G
23	BA	412	A
23	BA	444	C
23	BA	457	A
23	BA	470	A
23	BA	473	G
23	BA	480	A
23	BA	481	G
23	BA	482	A
23	BA	505	A
23	BA	508	G
23	BA	509	C
23	BA	530	G
23	BA	531	C
23	BA	532	A
23	BA	556	G
23	BA	563	G
23	BA	569	U
23	BA	573	G
23	BA	575	A
23	BA	598	G
23	BA	603	A
23	BA	615	G
23	BA	617	G
23	BA	620	G
23	BA	627	A

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
23	BA	637	A
23	BA	645	C
23	BA	646	A
23	BA	654	U
23	BA	655	A
23	BA	686	G
23	BA	695	G
23	BA	730	C
23	BA	740	U
23	BA	746	A
23	BA	747	U
23	BA	749	C
23	BA	764	A
23	BA	776	G
23	BA	779	U
23	BA	782	A
23	BA	784	A
23	BA	785	G
23	BA	790	C
23	BA	792	G
23	BA	800	A
23	BA	805	G
23	BA	812	C
23	BA	819	A
23	BA	827	U
23	BA	828	U
23	BA	846	C
23	BA	847	U
23	BA	859	G
23	BA	887	A
23	BA	889	C
23	BA	890	A
23	BA	896	A
23	BA	897	C
23	BA	910	A
23	BA	917	A
23	BA	931	G
23	BA	932	G
23	BA	941	A
23	BA	945	A
23	BA	946	G
23	BA	948	G

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
23	BA	959	A
23	BA	961	C
23	BA	973	A
23	BA	974(A)	G
23	BA	974(B)	C
23	BA	975	G
23	BA	983	A
23	BA	989	G
23	BA	990	A
23	BA	996	A
23	BA	1011	G
23	BA	1012	U
23	BA	1013	C
23	BA	1022	G
23	BA	1023	U
23	BA	1025	G
23	BA	1026	U
23	BA	1033	U
23	BA	1047	G
23	BA	1112	G
23	BA	1126	A
23	BA	1129	A
23	BA	1130	U
23	BA	1135	C
23	BA	1136	G
23	BA	1139	G
23	BA	114(B)	A
23	BA	1144	G
23	BA	1155	A
23	BA	1164	G
23	BA	1174	A
23	BA	1175	U
23	BA	1205	U
23	BA	1210	A
23	BA	1211	U
23	BA	1220	A
23	BA	1221	C
23	BA	1227	G
23	BA	1241	A
23	BA	1242	A
23	BA	1249	U
23	BA	1250	G

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
23	BA	1253	A
23	BA	1256	G
23	BA	1271	G
23	BA	1272	A
23	BA	1300	U
23	BA	1301	A
23	BA	1309	G
23	BA	1311	G
23	BA	1312	U
23	BA	1314	C
23	BA	1329	U
23	BA	1343	G
23	BA	1345	C
23	BA	1349	A
23	BA	1352	U
23	BA	1359	A
23	BA	1360	A
23	BA	1365	A
23	BA	1368	G
23	BA	1378	A
23	BA	1379	A
23	BA	1380	G
23	BA	1384	A
23	BA	1385	G
23	BA	1386	C
23	BA	1396	U
23	BA	1416	G
23	BA	1417	C
23	BA	1420	U
23	BA	1427	A
23	BA	1428	C
23	BA	144(B)	A
23	BA	1453	A
23	BA	1467	C
23	BA	1483	G
23	BA	1490	A
23	BA	1493	C
23	BA	1494	A
23	BA	1495	A
23	BA	1497	U
23	BA	1505	C
23	BA	1509	A

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
23	BA	1510	A
23	BA	1542	G
23	BA	1543	A
23	BA	1544	C
23	BA	1545	A
23	BA	1558	A
23	BA	1559	G
23	BA	1569	A
23	BA	1579	A
23	BA	1585	C
23	BA	1586	A
23	BA	1598	C
23	BA	1603	A
23	BA	1604	C
23	BA	1609	A
23	BA	1610	A
23	BA	1617	C
23	BA	1618	A
23	BA	1640	C
23	BA	1647	G
23	BA	1648	C
23	BA	1673	U
23	BA	1674	G
23	BA	1694	C
23	BA	1695	G
23	BA	1696	G
23	BA	1729	A
23	BA	1763	G
23	BA	1764	G
23	BA	1773	A
23	BA	1776	G
23	BA	1786	A
23	BA	1791	A
23	BA	1800	C
23	BA	1801	G
23	BA	1803	A
23	BA	1811	G
23	BA	1816	G
23	BA	1829	A
23	BA	1838	C
23	BA	1839	G
23	BA	1847	A

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
23	BA	1888	G
23	BA	1889	A
23	BA	1902	C
23	BA	1903	G
23	BA	1906	G
23	BA	1913	A
23	BA	1914	C
23	BA	1929	G
23	BA	1930	G
23	BA	1936	A
23	BA	1938	A
23	BA	1939	U
23	BA	1955	U
23	BA	1963	U
23	BA	1966	A
23	BA	1967	C
23	BA	1970	A
23	BA	1971	A
23	BA	1972	A
23	BA	1982	C
23	BA	1991	U
23	BA	1992	G
23	BA	1993	U
23	BA	1997	G
23	BA	2004	G
23	BA	2023	G
23	BA	2031	A
23	BA	2032	G
23	BA	2033	A
23	BA	2034	U
23	BA	2036	C
23	BA	2043	C
23	BA	2049	G
23	BA	2052	G
23	BA	2055	C
23	BA	2056	G
23	BA	2060	A
23	BA	2061	G
23	BA	2062	A
23	BA	2063	C
23	BA	2069	G
23	BA	2080	G

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
23	BA	2099	U
23	BA	2184	G
23	BA	2189	U
23	BA	2190	G
23	BA	2198	A
23	BA	2211	G
23	BA	2212	A
23	BA	2215	G
23	BA	2225	A
23	BA	2226	C
23	BA	2238	G
23	BA	2239	G
23	BA	2251	G
23	BA	2273	A
23	BA	2275	C
23	BA	2283	C
23	BA	2287	A
23	BA	2288	A
23	BA	2305	A
23	BA	2306	C
23	BA	2307	G
23	BA	2310	A
23	BA	2311	A
23	BA	2312	U
23	BA	2319	G
23	BA	2320	A
23	BA	2322	A
23	BA	2325	G
23	BA	2334	G
23	BA	2336	A
23	BA	2343	C
23	BA	2345	G
23	BA	2346	A
23	BA	2347	C
23	BA	2360	A
23	BA	2372	G
23	BA	2379	G
23	BA	2383	G
23	BA	2385	C
23	BA	2402	C
23	BA	2413	G
23	BA	2425	A

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
23	BA	2427	C
23	BA	2428	G
23	BA	2429	G
23	BA	2430	A
23	BA	2431	U
23	BA	2434	A
23	BA	2435	A
23	BA	2439	A
23	BA	2440	C
23	BA	2441	C
23	BA	2445	G
23	BA	2448	A
23	BA	2469	A
23	BA	2476	A
23	BA	2478	A
23	BA	2482	G
23	BA	2487	G
23	BA	2491	U
23	BA	2502	G
23	BA	2505	G
23	BA	2513	G
23	BA	2518	A
23	BA	2520	C
23	BA	2529	G
23	BA	2542	A
23	BA	2543	G
23	BA	2549	G
23	BA	2554	U
23	BA	2562	U
23	BA	2566	A
23	BA	2567	G
23	BA	2573	C
23	BA	2602	A
23	BA	2609	U
23	BA	2611	U
23	BA	2612	C
23	BA	2621	A
23	BA	2665	A
23	BA	2690	C
23	BA	2691	C
23	BA	2702	U
23	BA	2712	U

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
23	BA	712(B)	A
23	BA	2713	A
23	BA	2714	G
23	BA	2726	U
23	BA	2733	A
23	BA	2757	A
23	BA	2758	A
23	BA	2764	A
23	BA	2765	A
23	BA	2778	A
23	BA	2779	U
23	BA	2781	A
23	BA	2790	A
23	BA	2791	C
23	BA	2797	U
23	BA	2808	U
23	BA	2818	G
23	BA	2820	A
23	BA	2821	A
23	BA	2834	G
23	BA	2836	U
23	BA	2850	A
23	BA	2851	A
23	BA	2872	G
23	BA	2876	G
23	BA	2892	A
23	BA	2894	G
24	BB	12	C
24	BB	13	A
24	BB	15	A
24	BB	16	G
24	BB	42	C
24	BB	44	G
24	BB	45	A
24	BB	67	G
24	BB	73	A
24	BB	77	U
24	BB	88	C
24	BB	89(A)	G
24	BB	90	C
24	BB	108	C
24	BB	109	G

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Mol	Chain	Res	Type
24	BB	112	G
1	CA	9	G
1	CA	22	G
1	CA	32	A
1	CA	39	G
1	CA	41	G
1	CA	47	C
1	CA	48	C
1	CA	51	A
1	CA	61	G
1	CA	88	C
1	CA	99	C
1	CA	101	A
1	CA	108	G
1	CA	116	A
1	CA	120	A
1	CA	121	C
1	CA	131	C
1	CA	169	C
1	CA	182	U
1	CA	195	A
1	CA	197	A
1	CA	200	G
1	CA	209	U
1	CA	210	U
1	CA	243	A
1	CA	244	U
1	CA	247	G
1	CA	251	G
1	CA	266	G
1	CA	267	C
1	CA	289	G
1	CA	321	A
1	CA	328	C
1	CA	329	A
1	CA	332	G
1	CA	344	A
1	CA	345	C
1	CA	346	G
1	CA	352	C
1	CA	353	A
1	CA	354	G

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
1	CA	358	U
1	CA	367	U
1	CA	372	C
1	CA	373	A
1	CA	382	A
1	CA	387	U
1	CA	397	A
1	CA	398	C
1	CA	412	A
1	CA	413	G
1	CA	414	A
1	CA	422	C
1	CA	423	G
1	CA	429	U
1	CA	430	A
1	CA	439	A
1	CA	465	A
1	CA	467	G
1	CA	485	G
1	CA	496	A
1	CA	497	U
1	CA	511	C
1	CA	512	U
1	CA	518	C
1	CA	527	G
1	CA	531	U
1	CA	532	A
1	CA	533	A
1	CA	547	A
1	CA	559	A
1	CA	561	U
1	CA	562	C
1	CA	572	A
1	CA	573	A
1	CA	576	G
1	CA	596	C
1	CA	653	A
1	CA	687	A
1	CA	688	G
1	CA	703	G
1	CA	731	G
1	CA	755	G

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Mol	Chain	Res	Type
1	CA	777	A
1	CA	793	U
1	CA	794	A
1	CA	796	C
1	CA	816	A
1	CA	817	C
1	CA	818	G
1	CA	819	A
1	CA	828	A
1	CA	829	G
1	CA	833	U
1	CA	841	U
1	CA	842	C
1	CA	843	U
1	CA	855	G
1	CA	859	A
1	CA	873	A
1	CA	902	G
1	CA	914	A
1	CA	926	G
1	CA	927	G
1	CA	934	C
1	CA	935	A
1	CA	960	U
1	CA	961	U
1	CA	969	A
1	CA	971	G
1	CA	974	A
1	CA	976	G
1	CA	977	A
1	CA	980	C
1	CA	981	U
1	CA	992	U
1	CA	993	G
1	CA	1004	A
1	CA	1006	C
1	CA	1045	C
1	CA	1054	C
1	CA	1065	U
1	CA	1066	C
1	CA	1068	G
1	CA	1081	G

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Mol	Chain	Res	Type
1	CA	1094	G
1	CA	1095	U
1	CA	1101	A
1	CA	1113	C
1	CA	1117	G
1	CA	1118	C
1	CA	1125	U
1	CA	1126	U
1	CA	1129	C
1	CA	1130	A
1	CA	1131	G
1	CA	1136	U
1	CA	1137	C
1	CA	1139	G
1	CA	1146	A
1	CA	1151	A
1	CA	1152	A
1	CA	1159	U
1	CA	1160	G
1	CA	1181	G
1	CA	1182	G
1	CA	1193	G
1	CA	1196	U
1	CA	1197	G
1	CA	1202	G
1	CA	1212	U
1	CA	1213	A
1	CA	1214	C
1	CA	1224	G
1	CA	1225	A
1	CA	1238	A
1	CA	1240	U
1	CA	1241	G
1	CA	1256	A
1	CA	1257	U
1	CA	1280	A
1	CA	1281	U
1	CA	1282	C
1	CA	1285	A
1	CA	1286	A
1	CA	1287	A
1	CA	1300	G

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Mol	Chain	Res	Type
1	CA	1301	U
1	CA	1302	U
1	CA	1305	G
1	CA	1317	C
1	CA	1320	C
1	CA	1322	C
1	CA	1331	G
1	CA	1335	C
1	CA	1346	A
1	CA	1347	G
1	CA	136(B)	C
1	CA	1363	A
1	CA	1364	U
1	CA	1400	C
1	CA	1401	G
1	CA	1419	G
1	CA	1442	G
1	CA	1443	G
1	CA	1446	A
1	CA	1447	G
1	CA	1451	A
1	CA	1452	C
1	CA	1454	G
1	CA	1493	A
1	CA	1494	G
1	CA	1497	G
1	CA	1499	A
1	CA	1502	A
1	CA	1503	A
1	CA	1504	G
1	CA	1505	G
1	CA	1506	U
1	CA	1517	G
1	CA	1519	A
1	CA	1520	G
1	CA	1528	U
1	CA	1529	G
1	CA	1530	G
22	CV	6172	U
22	CV	6189	A
23	DA	10	G
23	DA	35	G

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
23	DA	46	C
23	DA	64	A
23	DA	71	A
23	DA	72	U
23	DA	74	A
23	DA	75	G
23	DA	84	A
23	DA	88	G
23	DA	102	G
23	DA	118	A
23	DA	119	A
23	DA	120	U
23	DA	125	G
23	DA	131	G
23	DA	138	G
23	DA	140	A
23	DA	196	A
23	DA	197	A
23	DA	199	A
23	DA	205	G
23	DA	215	G
23	DA	216	A
23	DA	221	A
23	DA	222	A
23	DA	228	A
23	DA	229	A
23	DA	230	U
23	DA	245	G
23	DA	248	G
23	DA	252	G
23	DA	269	U
23	DA	270(K)	G
23	DA	270(M)	U
23	DA	270(N)	U
23	DA	270(O)	G
23	DA	270(R)	C
23	DA	271(D)	U
23	DA	271	G
23	DA	274	G
23	DA	275	G
23	DA	276	A
23	DA	277	C

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
23	DA	278	A
23	DA	279	C
23	DA	283	A
23	DA	302	C
23	DA	311	A
23	DA	323	G
23	DA	324	A
23	DA	329	G
23	DA	330	A
23	DA	332	A
23	DA	333	G
23	DA	343	C
23	DA	352	G
23	DA	353	G
23	DA	386	G
23	DA	396	G
23	DA	405	U
23	DA	411	G
23	DA	412	A
23	DA	444	C
23	DA	457	A
23	DA	470	A
23	DA	480	A
23	DA	481	G
23	DA	482	A
23	DA	505	A
23	DA	508	G
23	DA	509	C
23	DA	530	G
23	DA	531	C
23	DA	532	A
23	DA	556	G
23	DA	563	G
23	DA	569	U
23	DA	573	G
23	DA	575	A
23	DA	598	G
23	DA	603	A
23	DA	615	G
23	DA	617	G
23	DA	620	G
23	DA	627	A

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
23	DA	637	A
23	DA	645	C
23	DA	646	A
23	DA	654	U
23	DA	655	A
23	DA	686	G
23	DA	695	G
23	DA	730	C
23	DA	746	A
23	DA	747	U
23	DA	749	C
23	DA	764	A
23	DA	776	G
23	DA	777	A
23	DA	779	U
23	DA	782	A
23	DA	784	A
23	DA	785	G
23	DA	790	C
23	DA	792	G
23	DA	800	A
23	DA	805	G
23	DA	812	C
23	DA	819	A
23	DA	827	U
23	DA	828	U
23	DA	846	C
23	DA	847	U
23	DA	859	G
23	DA	887	A
23	DA	889	C
23	DA	890	A
23	DA	896	A
23	DA	897	C
23	DA	910	A
23	DA	917	A
23	DA	931	G
23	DA	932	G
23	DA	941	A
23	DA	945	A
23	DA	946	G
23	DA	948	G

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Mol	Chain	Res	Type
23	DA	959	A
23	DA	961	C
23	DA	973	A
23	DA	974(A)	G
23	DA	974(B)	C
23	DA	975	G
23	DA	983	A
23	DA	989	G
23	DA	990	A
23	DA	996	A
23	DA	1011	G
23	DA	1012	U
23	DA	1013	C
23	DA	1022	G
23	DA	1023	U
23	DA	1025	G
23	DA	1026	U
23	DA	1033	U
23	DA	1047	G
23	DA	1112	G
23	DA	1126	A
23	DA	1129	A
23	DA	1130	U
23	DA	1135	C
23	DA	1136	G
23	DA	1139	G
23	DA	114(B)	A
23	DA	1144	G
23	DA	1155	A
23	DA	1164	G
23	DA	1174	A
23	DA	1175	U
23	DA	1205	U
23	DA	1210	A
23	DA	1211	U
23	DA	1220	A
23	DA	1221	C
23	DA	1227	G
23	DA	1241	A
23	DA	1242	A
23	DA	1249	U
23	DA	1250	G

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
23	DA	1253	A
23	DA	1256	G
23	DA	1271	G
23	DA	1272	A
23	DA	1300	U
23	DA	1301	A
23	DA	1309	G
23	DA	1312	U
23	DA	1314	C
23	DA	1329	U
23	DA	1343	G
23	DA	1345	C
23	DA	1349	A
23	DA	1352	U
23	DA	1359	A
23	DA	1360	A
23	DA	1365	A
23	DA	1368	G
23	DA	1378	A
23	DA	1379	A
23	DA	1380	G
23	DA	1384	A
23	DA	1385	G
23	DA	1386	C
23	DA	1396	U
23	DA	1416	G
23	DA	1417	C
23	DA	1420	U
23	DA	1427	A
23	DA	1428	C
23	DA	144(B)	A
23	DA	1453	A
23	DA	1455	G
23	DA	1467	C
23	DA	1483	G
23	DA	1490	A
23	DA	1493	C
23	DA	1494	A
23	DA	1495	A
23	DA	1497	U
23	DA	1505	C
23	DA	1509	A

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
23	DA	1510	A
23	DA	1542	G
23	DA	1543	A
23	DA	1544	C
23	DA	1545	A
23	DA	1558	A
23	DA	1559	G
23	DA	1569	A
23	DA	1579	A
23	DA	1585	C
23	DA	1586	A
23	DA	1598	C
23	DA	1603	A
23	DA	1604	C
23	DA	1609	A
23	DA	1610	A
23	DA	1617	C
23	DA	1618	A
23	DA	1640	C
23	DA	1647	G
23	DA	1648	C
23	DA	1673	U
23	DA	1674	G
23	DA	1694	C
23	DA	1695	G
23	DA	1696	G
23	DA	1729	A
23	DA	1763	G
23	DA	1764	G
23	DA	1773	A
23	DA	1776	G
23	DA	1786	A
23	DA	1791	A
23	DA	1800	C
23	DA	1801	G
23	DA	1803	A
23	DA	1811	G
23	DA	1816	G
23	DA	1829	A
23	DA	1838	C
23	DA	1839	G
23	DA	1847	A

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
23	DA	1888	G
23	DA	1889	A
23	DA	1902	C
23	DA	1903	G
23	DA	1906	G
23	DA	1913	A
23	DA	1914	C
23	DA	1929	G
23	DA	1930	G
23	DA	1936	A
23	DA	1938	A
23	DA	1939	U
23	DA	1955	U
23	DA	1963	U
23	DA	1966	A
23	DA	1967	C
23	DA	1970	A
23	DA	1971	A
23	DA	1972	A
23	DA	1982	C
23	DA	1991	U
23	DA	1992	G
23	DA	1993	U
23	DA	1997	G
23	DA	2004	G
23	DA	2023	G
23	DA	2031	A
23	DA	2032	G
23	DA	2033	A
23	DA	2034	U
23	DA	2036	C
23	DA	2043	C
23	DA	2049	G
23	DA	2052	G
23	DA	2055	C
23	DA	2056	G
23	DA	2060	A
23	DA	2061	G
23	DA	2062	A
23	DA	2063	C
23	DA	2069	G
23	DA	2080	G

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
23	DA	2099	U
23	DA	2184	G
23	DA	2189	U
23	DA	2190	G
23	DA	2198	A
23	DA	2211	G
23	DA	2212	A
23	DA	2215	G
23	DA	2225	A
23	DA	2226	C
23	DA	2238	G
23	DA	2239	G
23	DA	2251	G
23	DA	2273	A
23	DA	2275	C
23	DA	2283	C
23	DA	2287	A
23	DA	2288	A
23	DA	2305	A
23	DA	2306	C
23	DA	2307	G
23	DA	2310	A
23	DA	2311	A
23	DA	2312	U
23	DA	2319	G
23	DA	2320	A
23	DA	2322	A
23	DA	2325	G
23	DA	2334	G
23	DA	2336	A
23	DA	2343	C
23	DA	2345	G
23	DA	2346	A
23	DA	2347	C
23	DA	2360	A
23	DA	2372	G
23	DA	2379	G
23	DA	2383	G
23	DA	2385	C
23	DA	2402	C
23	DA	2413	G
23	DA	2425	A

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
23	DA	2427	C
23	DA	2428	G
23	DA	2429	G
23	DA	2430	A
23	DA	2431	U
23	DA	2434	A
23	DA	2435	A
23	DA	2439	A
23	DA	2440	C
23	DA	2441	C
23	DA	2445	G
23	DA	2448	A
23	DA	2469	A
23	DA	2476	A
23	DA	2478	A
23	DA	2482	G
23	DA	2487	G
23	DA	2491	U
23	DA	2502	G
23	DA	2505	G
23	DA	2513	G
23	DA	2518	A
23	DA	2520	C
23	DA	2529	G
23	DA	2542	A
23	DA	2543	G
23	DA	2554	U
23	DA	2562	U
23	DA	2566	A
23	DA	2567	G
23	DA	2572	A
23	DA	2573	C
23	DA	2602	A
23	DA	2609	U
23	DA	2611	U
23	DA	2612	C
23	DA	2621	A
23	DA	2665	A
23	DA	2690	C
23	DA	2691	C
23	DA	2702	U
23	DA	2712	U

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Mol	Chain	Res	Type
23	DA	712(B)	A
23	DA	2713	A
23	DA	2714	G
23	DA	2726	U
23	DA	2733	A
23	DA	2757	A
23	DA	2758	A
23	DA	2764	A
23	DA	2765	A
23	DA	2778	A
23	DA	2779	U
23	DA	2781	A
23	DA	2790	A
23	DA	2791	C
23	DA	2797	U
23	DA	2808	U
23	DA	2818	G
23	DA	2820	A
23	DA	2821	A
23	DA	2834	G
23	DA	2836	U
23	DA	2850	A
23	DA	2851	A
23	DA	2872	G
23	DA	2876	G
23	DA	2892	A
23	DA	2894	G
24	DB	12	C
24	DB	13	A
24	DB	15	A
24	DB	16	G
24	DB	42	C
24	DB	44	G
24	DB	45	A
24	DB	67	G
24	DB	73	A
24	DB	77	U
24	DB	88	C
24	DB	89(A)	G
24	DB	90	C
24	DB	108	C
24	DB	109	G

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Mol	Chain	Res	Type
24	DB	112	G

All (215) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
1	AA	7	G
1	AA	99	C
1	AA	119	A
1	AA	121	C
1	AA	243	A
1	AA	250	A
1	AA	266	G
1	AA	328	C
1	AA	353	A
1	AA	366	C
1	AA	428	G
1	AA	429	U
1	AA	438	G
1	AA	496	A
1	AA	560	U
1	AA	561	U
1	AA	562	C
1	AA	687	A
1	AA	913	A
1	AA	1064	G
1	AA	1065	U
1	AA	1067	A
1	AA	1129	C
1	AA	1145	C
1	AA	1201	A
1	AA	1281	U
1	AA	1285	A
1	AA	1300	G
1	AA	1400	C
1	AA	1492	A
1	AA	1493	A
1	AA	1498	U
1	AA	1504	G
1	AA	1528	U
1	AA	1529	G
22	AV	6171	U
22	AV	6188	C

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
23	BA	102	G
23	BA	119	A
23	BA	120	U
23	BA	131	G
23	BA	199	A
23	BA	221	A
23	BA	278	A
23	BA	331	A
23	BA	332	A
23	BA	479	A
23	BA	481	G
23	BA	503	A
23	BA	616	A
23	BA	682	G
23	BA	685	A
23	BA	746	A
23	BA	764	A
23	BA	776	G
23	BA	846	C
23	BA	858	U
23	BA	859	G
23	BA	933	A
23	BA	945	A
23	BA	948	G
23	BA	961	C
23	BA	974(A)	G
23	BA	989	G
23	BA	1022	G
23	BA	1157	G
23	BA	1210	A
23	BA	1221	C
23	BA	1253	A
23	BA	1311	G
23	BA	1314	C
23	BA	1343	G
23	BA	1378	A
23	BA	1379	A
23	BA	1427	A
23	BA	1494	A
23	BA	1542	G
23	BA	1558	A
23	BA	1579	A

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
23	BA	1603	A
23	BA	1608	A
23	BA	1694	C
23	BA	1786	A
23	BA	1800	C
23	BA	1929	G
23	BA	1930	G
23	BA	1937	A
23	BA	2033	A
23	BA	2051	A
23	BA	2062	A
23	BA	2098	U
23	BA	2225	A
23	BA	2250	G
23	BA	2311	A
23	BA	2426	A
23	BA	2427	C
23	BA	2439	A
23	BA	2468	G
23	BA	2481	G
23	BA	2581	G
23	BA	2585	U
23	BA	2689	U
23	BA	2756	U
23	BA	2849	U
23	BA	2850	A
23	BA	2873	A
24	BB	66	A
1	CA	7	G
1	CA	99	C
1	CA	119	A
1	CA	121	C
1	CA	243	A
1	CA	250	A
1	CA	266	G
1	CA	328	C
1	CA	353	A
1	CA	366	C
1	CA	428	G
1	CA	429	U
1	CA	438	G
1	CA	496	A

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Mol	Chain	Res	Type
1	CA	560	U
1	CA	561	U
1	CA	562	C
1	CA	687	A
1	CA	913	A
1	CA	1064	G
1	CA	1065	U
1	CA	1067	A
1	CA	1129	C
1	CA	1145	C
1	CA	1201	A
1	CA	1281	U
1	CA	1285	A
1	CA	1300	G
1	CA	1400	C
1	CA	1492	A
1	CA	1493	A
1	CA	1498	U
1	CA	1504	G
1	CA	1528	U
1	CA	1529	G
22	CV	6171	U
22	CV	6188	C
23	DA	102	G
23	DA	119	A
23	DA	120	U
23	DA	131	G
23	DA	199	A
23	DA	221	A
23	DA	278	A
23	DA	331	A
23	DA	332	A
23	DA	343	C
23	DA	479	A
23	DA	481	G
23	DA	503	A
23	DA	616	A
23	DA	682	G
23	DA	685	A
23	DA	746	A
23	DA	764	A
23	DA	776	G

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Mol	Chain	Res	Type
23	DA	846	C
23	DA	858	U
23	DA	859	G
23	DA	933	A
23	DA	945	A
23	DA	948	G
23	DA	961	C
23	DA	974(A)	G
23	DA	989	G
23	DA	1022	G
23	DA	1157	G
23	DA	1210	A
23	DA	1221	C
23	DA	1253	A
23	DA	1311	G
23	DA	1314	C
23	DA	1343	G
23	DA	1378	A
23	DA	1379	A
23	DA	1427	A
23	DA	1494	A
23	DA	1542	G
23	DA	1558	A
23	DA	1579	A
23	DA	1603	A
23	DA	1608	A
23	DA	1694	C
23	DA	1800	C
23	DA	1929	G
23	DA	1930	G
23	DA	1937	A
23	DA	1980	G
23	DA	2033	A
23	DA	2051	A
23	DA	2062	A
23	DA	2098	U
23	DA	2225	A
23	DA	2250	G
23	DA	2311	A
23	DA	2426	A
23	DA	2427	C
23	DA	2439	A

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Mol	Chain	Res	Type
23	DA	2468	G
23	DA	2481	G
23	DA	2581	G
23	DA	2585	U
23	DA	2689	U
23	DA	2756	U
23	DA	2849	U
23	DA	2850	A
23	DA	2873	A
24	DB	66	A

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

There are no non-standard protein/DNA/RNA residues in this entry.

## 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

Of 2661 ligands modelled in this entry, 2661 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 ⓘ

There are no such residues in this entry.

## 5.8 Polymer linkage issues ⓘ

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, 95<sup>th</sup> 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(Å <sup>2</sup> )	Q<0.9
1	AA	1506/1506 (100%)	0.10	58 (3%)	43	35	66, 119, 257, 429	0
1	CA	1506/1506 (100%)	0.32	99 (6%)	22	16	75, 164, 286, 403	0
2	AB	234/234 (100%)	1.53	80 (34%)	0	0	88, 178, 260, 354	0
2	CB	234/234 (100%)	1.37	73 (31%)	1	1	106, 202, 279, 372	0
3	AC	206/206 (100%)	0.59	22 (10%)	8	7	83, 165, 233, 293	0
3	CC	206/206 (100%)	0.67	32 (15%)	3	3	124, 198, 272, 337	0
4	AD	208/208 (100%)	0.04	4 (1%)	70	60	45, 119, 180, 228	0
4	CD	208/208 (100%)	0.48	20 (9%)	10	9	96, 186, 253, 314	0
5	AE	151/151 (100%)	0.00	3 (1%)	68	59	71, 118, 191, 293	0
5	CE	151/151 (100%)	0.20	6 (3%)	42	33	72, 165, 223, 282	0
6	AF	101/101 (100%)	0.72	19 (18%)	2	2	87, 160, 207, 278	0
6	CF	101/101 (100%)	0.03	5 (4%)	32	25	60, 130, 195, 271	0
7	AG	155/155 (100%)	0.98	30 (19%)	1	2	134, 201, 260, 328	0
7	CG	155/155 (100%)	1.10	34 (21%)	1	1	124, 213, 271, 320	0
8	AH	138/138 (100%)	0.46	15 (10%)	7	7	60, 122, 190, 265	0
8	CH	138/138 (100%)	0.70	19 (13%)	4	4	101, 171, 217, 257	0
9	AI	127/127 (100%)	1.82	49 (38%)	0	0	119, 228, 303, 333	0
9	CI	127/127 (100%)	1.48	39 (30%)	1	1	148, 229, 297, 370	0
10	AJ	98/98 (100%)	2.01	46 (46%)	0	0	113, 203, 294, 334	0
10	CJ	98/98 (100%)	1.92	40 (40%)	0	0	133, 232, 290, 339	0
11	AK	119/119 (100%)	-0.07	9 (7%)	17	14	66, 136, 219, 274	0
11	CK	119/119 (100%)	0.27	8 (6%)	21	16	87, 145, 207, 287	0
12	AL	124/124 (100%)	0.60	9 (7%)	18	14	56, 108, 182, 328	0
12	CL	124/124 (100%)	0.48	12 (9%)	10	9	77, 139, 236, 271	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
13	AM	116/116 (100%)	1.07	25 (21%) 1 1	129, 216, 277, 325	0
13	CM	116/116 (100%)	2.08	51 (43%) 0 0	119, 238, 307, 372	0
14	AN	60/60 (100%)	1.46	17 (28%) 1 1	94, 178, 243, 278	0
14	CN	60/60 (100%)	1.91	25 (41%) 0 0	110, 207, 261, 344	0
15	AO	88/88 (100%)	-0.06	0 100 100	73, 127, 187, 281	0
15	CO	88/88 (100%)	0.12	5 (5%) 27 21	73, 138, 201, 244	0
16	AP	83/83 (100%)	0.31	4 (4%) 34 27	67, 109, 175, 245	0
16	CP	83/83 (100%)	0.87	23 (27%) 1 1	128, 171, 247, 277	0
17	AQ	99/99 (100%)	-0.12	0 100 100	63, 112, 169, 222	0
17	CQ	99/99 (100%)	0.70	15 (15%) 3 3	99, 155, 196, 228	0
18	AR	70/70 (100%)	1.24	18 (25%) 1 1	82, 149, 241, 325	0
18	CR	70/70 (100%)	0.82	10 (14%) 4 4	76, 145, 204, 232	0
19	AS	78/78 (100%)	1.92	32 (41%) 0 0	133, 212, 274, 295	0
19	CS	78/78 (100%)	2.12	35 (44%) 0 0	173, 235, 302, 347	0
20	AT	99/99 (100%)	0.42	8 (8%) 15 12	65, 131, 220, 241	0
20	CT	99/99 (100%)	2.09	47 (47%) 0 0	103, 178, 284, 301	0
21	AU	24/24 (100%)	2.38	13 (54%) 0 0	122, 220, 292, 333	0
21	CU	24/24 (100%)	2.01	9 (37%) 0 0	173, 236, 317, 346	0
22	AV	35/35 (100%)	1.52	12 (34%) 0 0	96, 223, 325, 366	0
22	CV	35/35 (100%)	1.13	6 (17%) 2 2	117, 192, 362, 386	0
23	BA	2760/2879 (95%)	-0.04	67 (2%) 62 52	38, 86, 205, 371	0
23	DA	2760/2879 (95%)	-0.10	44 (1%) 74 65	33, 77, 196, 377	0
24	BB	119/119 (100%)	-0.02	0 100 100	83, 149, 196, 251	0
24	DB	119/119 (100%)	0.03	6 (5%) 32 25	87, 132, 198, 274	0
25	BC	271/271 (100%)	-0.05	4 (1%) 76 67	24, 81, 150, 209	0
25	DC	271/271 (100%)	-0.30	0 100 100	26, 72, 144, 220	0
26	BD	204/204 (100%)	0.14	3 (1%) 76 67	45, 96, 173, 331	0
26	DD	204/204 (100%)	-0.02	3 (1%) 76 67	36, 90, 189, 305	0
27	BE	202/202 (100%)	0.12	4 (1%) 68 59	32, 94, 172, 311	0
27	DE	202/202 (100%)	0.18	8 (3%) 42 33	20, 85, 182, 249	0
28	BF	181/181 (100%)	0.50	19 (10%) 8 8	107, 202, 262, 297	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
28	DF	181/181 (100%)	0.48	20 (11%) 7 7	109, 198, 281, 324	0
29	BG	159/159 (100%)	1.28	47 (29%) 1 1	102, 192, 269, 331	0
29	DG	159/159 (100%)	0.20	10 (6%) 23 18	61, 117, 190, 269	0
30	BH	145/145 (100%)	1.78	42 (28%) 1 1	60, 224, 376, 453	0
30	DH	145/145 (100%)	1.73	49 (33%) 0 0	60, 210, 372, 482	0
31	BI	32/65 (49%)	6.39	31 (96%) 0 0	163, 256, 325, 352	0
31	DI	32/65 (49%)	2.47	21 (65%) 0 0	176, 235, 294, 325	0
32	BJ	137/137 (100%)	0.30	3 (2%) 65 55	58, 105, 182, 221	0
32	DJ	137/137 (100%)	0.03	1 (0%) 89 82	55, 101, 170, 219	0
33	BK	122/122 (100%)	-0.22	0 100 100	48, 90, 142, 194	0
33	DK	122/122 (100%)	-0.21	0 100 100	41, 90, 153, 251	0
34	BL	146/146 (100%)	0.50	10 (6%) 20 16	40, 114, 204, 306	0
34	DL	146/146 (100%)	0.42	11 (7%) 17 14	27, 111, 207, 324	0
35	BM	136/136 (100%)	0.12	4 (2%) 55 45	57, 110, 205, 344	0
35	DM	136/136 (100%)	0.41	8 (5%) 26 20	46, 106, 199, 388	0
36	BN	117/117 (100%)	0.14	2 (1%) 73 64	44, 92, 166, 282	0
36	DN	117/117 (100%)	0.09	2 (1%) 73 64	41, 89, 173, 285	0
37	BO	98/98 (100%)	2.05	46 (46%) 0 0	89, 155, 230, 299	0
37	DO	98/98 (100%)	0.73	18 (18%) 2 2	80, 148, 221, 299	0
38	BP	137/137 (100%)	0.30	6 (4%) 38 29	52, 115, 216, 248	0
38	DP	137/137 (100%)	0.37	11 (8%) 15 12	60, 119, 249, 299	0
39	BQ	116/116 (100%)	-0.06	2 (1%) 73 64	43, 88, 163, 244	0
39	DQ	116/116 (100%)	-0.28	0 100 100	34, 84, 154, 205	0
40	BR	101/101 (100%)	-0.18	0 100 100	55, 134, 187, 294	0
40	DR	101/101 (100%)	-0.26	0 100 100	52, 132, 197, 321	0
41	BS	112/112 (100%)	0.48	5 (4%) 37 29	45, 78, 151, 250	0
41	DS	112/112 (100%)	0.03	1 (0%) 85 78	40, 72, 166, 277	0
42	BT	92/92 (100%)	0.09	1 (1%) 82 73	57, 107, 174, 204	0
42	DT	92/92 (100%)	0.10	0 100 100	41, 78, 161, 204	0
43	BU	100/100 (100%)	0.48	12 (12%) 6 6	59, 139, 289, 344	0
43	DU	100/100 (100%)	0.71	10 (10%) 9 9	49, 119, 264, 373	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
44	BV	188/188 (100%)	0.67	27 (14%) 3 4	80, 160, 224, 277	0
44	DV	188/188 (100%)	0.22	13 (6%) 20 16	67, 154, 220, 254	0
45	BW	76/76 (100%)	-0.07	3 (3%) 43 35	61, 101, 160, 258	0
45	DW	76/76 (100%)	0.20	2 (2%) 59 49	54, 101, 167, 254	0
46	BX	88/88 (100%)	0.06	3 (3%) 49 40	50, 106, 191, 346	0
46	DX	88/88 (100%)	-0.04	1 (1%) 82 73	31, 83, 196, 340	0
47	BY	62/62 (100%)	0.23	6 (9%) 10 9	61, 141, 244, 287	0
47	DY	62/62 (100%)	0.45	5 (8%) 15 12	45, 99, 224, 304	0
48	BZ	59/59 (100%)	1.41	13 (22%) 1 1	47, 102, 174, 342	0
48	DZ	59/59 (100%)	1.43	15 (25%) 1 1	45, 102, 189, 335	0
49	B1	30/30 (100%)	1.78	9 (30%) 1 1	161, 244, 282, 323	0
49	D1	30/30 (100%)	0.96	4 (13%) 4 5	187, 262, 306, 326	0
50	B2	52/52 (100%)	-0.29	1 (1%) 70 60	36, 93, 189, 273	0
50	D2	52/52 (100%)	-0.37	1 (1%) 70 60	24, 93, 214, 262	0
51	B3	44/44 (100%)	6.82	43 (97%) 0 0	151, 254, 304, 322	0
51	D3	44/44 (100%)	7.53	43 (97%) 0 0	191, 247, 298, 313	0
52	B4	48/48 (100%)	0.14	1 (2%) 67 58	43, 66, 132, 297	0
52	D4	48/48 (100%)	-0.27	0 100 100	23, 45, 122, 217	0
53	B5	63/63 (100%)	-0.20	1 (1%) 74 65	47, 94, 171, 222	0
53	D5	63/63 (100%)	0.18	0 100 100	43, 92, 170, 209	0
All	All	20232/20536 (98%)	0.37	1808 (8%) 12 10	20, 120, 258, 482	0

All (1808) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
43	DU	52	SER	23.5
1	AA	85	U	22.6
30	BH	84	GLY	21.6
30	DH	90	GLY	18.4
1	AA	82	U	17.7
1	AA	84	U	17.0
51	B3	49	HIS	16.9
51	B3	39	TYR	16.1
30	BH	85	GLU	15.8
35	BM	140	ALA	13.9

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Mol	Chain	Res	Type	RSRZ
51	D3	38	LYS	13.7
51	B3	40	CYS	13.6
51	B3	26	ASN	13.3
51	B3	13	CYS	12.9
51	D3	35	GLU	12.5
51	D3	20	ASN	12.4
30	DH	88	ILE	12.3
51	D3	49	HIS	12.3
1	AA	81	G	12.1
51	D3	12	GLU	11.8
51	D3	37	ARG	11.3
51	B3	47	THR	11.2
1	AA	86	U	11.2
51	D3	47	THR	11.1
30	BH	111	PRO	11.1
51	D3	21	TYR	11.1
31	BI	11	ALA	11.0
51	D3	13	CYS	10.8
51	D3	14	THR	10.8
51	B3	41	PRO	10.7
51	D3	39	TYR	10.4
51	D3	36	LEU	10.4
30	BH	86	THR	10.4
31	BI	8	GLU	10.3
51	B3	24	GLU	10.3
51	D3	22	ALA	10.0
30	DH	80	PRO	9.9
31	BI	66	LEU	9.8
51	B3	16	CYS	9.6
10	CJ	73	ASP	9.3
51	D3	41	PRO	9.2
20	CT	103	GLY	9.2
30	BH	119	PRO	9.2
31	BI	64	LYS	9.2
45	DW	85	ALA	9.1
51	D3	50	ARG	9.0
51	D3	43	CYS	8.9
31	BI	6	ASN	8.8
9	CI	8	GLY	8.6
17	CQ	8	GLY	8.6
47	DY	15	LYS	8.6
31	BI	63	LEU	8.6

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Mol	Chain	Res	Type	RSRZ
48	BZ	1	MET	8.5
51	D3	40	CYS	8.4
20	CT	102	GLY	8.4
51	B3	51	GLU	8.4
31	BI	61	LEU	8.3
31	BI	4	LYS	8.3
31	BI	3	ASN	8.3
19	CS	75	ALA	8.2
19	CS	41	VAL	8.2
2	CB	96	ARG	8.1
31	BI	5	ARG	8.1
3	CC	206	GLU	8.1
51	D3	52	VAL	8.1
51	B3	12	GLU	7.9
30	BH	121	LYS	7.9
2	CB	148	TYR	7.9
20	CT	60	GLU	7.8
51	B3	15	GLU	7.8
30	BH	68	LEU	7.7
9	CI	65	VAL	7.7
19	AS	4	SER	7.7
31	BI	9	LEU	7.6
31	BI	60	ARG	7.6
31	BI	15	GLU	7.6
51	D3	26	ASN	7.6
14	CN	12	ARG	7.6
51	D3	11	LEU	7.6
35	BM	141	GLN	7.5
51	B3	38	LYS	7.5
13	CM	8	GLU	7.5
22	AV	6150	U	7.5
35	DM	140	ALA	7.4
10	CJ	71	LEU	7.4
31	BI	12	THR	7.4
51	D3	44	ARG	7.4
35	DM	141	GLN	7.4
51	D3	18	ARG	7.3
30	BH	72	LEU	7.3
43	DU	53	PRO	7.3
51	B3	14	THR	7.3
13	CM	30	ALA	7.3
51	B3	17	LYS	7.3

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Mol	Chain	Res	Type	RSRZ
51	D3	24	GLU	7.3
31	BI	67	GLY	7.2
30	DH	94	ALA	7.2
10	CJ	35	SER	7.2
51	D3	19	ARG	7.1
43	DU	50	ARG	7.1
19	CS	74	PHE	7.1
14	AN	13	THR	7.1
48	DZ	1	MET	7.1
31	BI	65	GLU	7.0
2	CB	73	THR	7.0
23	DA	1535	U	7.0
43	DU	51	VAL	7.0
21	AU	23	PRO	6.9
2	CB	170	GLU	6.9
51	D3	48	VAL	6.9
19	CS	40	ILE	6.8
9	AI	91	ASP	6.8
51	B3	22	ALA	6.8
20	CT	51	GLU	6.8
23	BA	1046	A	6.8
29	BG	55	PRO	6.8
23	BA	2108	C	6.8
20	CT	59	ALA	6.8
31	BI	59	ILE	6.7
51	D3	15	GLU	6.7
2	CB	97	TRP	6.7
37	BO	58	LEU	6.7
10	AJ	27	ALA	6.7
51	B3	42	TRP	6.7
23	BA	2211	G	6.6
9	CI	64	THR	6.6
51	B3	25	LYS	6.6
19	AS	74	PHE	6.6
10	CJ	72	VAL	6.6
8	CH	132	GLU	6.6
18	AR	46	GLU	6.6
51	B3	46	HIS	6.6
51	B3	28	ARG	6.6
13	CM	15	VAL	6.6
10	AJ	73	ASP	6.6
30	DH	89	TYR	6.5

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
2	AB	165	VAL	6.4
31	BI	62	ALA	6.4
19	CS	48	THR	6.4
14	CN	13	THR	6.4
51	B3	50	ARG	6.4
9	AI	64	THR	6.4
51	D3	16	CYS	6.3
7	CG	77	SER	6.3
2	CB	188	ALA	6.3
23	DA	2801	A	6.3
30	BH	100	ALA	6.3
51	D3	46	HIS	6.3
51	D3	32	ASN	6.3
30	DH	132	PRO	6.2
19	AS	81	ARG	6.2
9	CI	7	THR	6.2
23	DA	6	A	6.2
13	CM	24	GLY	6.2
51	D3	51	GLU	6.2
30	DH	120	ILE	6.1
31	DI	62	ALA	6.1
37	BO	52	SER	6.1
2	CB	101	MET	6.1
20	AT	64	ASP	6.1
23	BA	2801	A	6.1
31	BI	14	LYS	6.1
19	CS	71	LEU	6.1
22	AV	6188	C	6.0
20	CT	55	ILE	6.0
6	AF	89	MET	6.0
3	CC	207	VAL	6.0
51	B3	44	ARG	6.0
21	CU	5	ASP	6.0
51	B3	20	ASN	5.9
23	DA	1174	A	5.9
1	AA	1129	C	5.9
19	AS	49	ILE	5.9
44	BV	186	GLU	5.9
51	B3	32	ASN	5.9
20	CT	100	ILE	5.9
7	CG	81	GLY	5.9
23	BA	2892	A	5.9

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Mol	Chain	Res	Type	RSRZ
37	BO	87	PHE	5.9
10	AJ	72	VAL	5.8
20	CT	104	LEU	5.8
7	CG	82	GLY	5.8
30	DH	107	ILE	5.8
51	D3	9	LEU	5.8
23	BA	2799	A	5.8
30	DH	92	VAL	5.7
10	CJ	38	ILE	5.7
51	D3	34	LEU	5.7
30	DH	122	GLU	5.7
1	AA	1000	A	5.7
1	CA	84	U	5.7
21	CU	17	THR	5.7
31	BI	7	VAL	5.7
9	AI	82	ALA	5.7
20	CT	52	ALA	5.7
9	CI	29	ASN	5.7
9	CI	127	LYS	5.7
31	BI	16	ASN	5.7
13	CM	5	ALA	5.6
31	DI	63	LEU	5.6
51	D3	23	THR	5.6
31	DI	66	LEU	5.6
51	B3	48	VAL	5.6
9	AI	62	TYR	5.6
23	DA	2107	C	5.6
30	BH	79	ILE	5.6
24	DB	52	A	5.6
1	AA	87	A	5.6
51	B3	21	TYR	5.6
13	CM	117	VAL	5.5
12	CL	71	GLY	5.5
19	AS	75	ALA	5.5
9	AI	85	LEU	5.5
23	DA	888	C	5.5
13	CM	65	LYS	5.5
29	BG	57	ASP	5.5
9	AI	17	VAL	5.5
2	CB	28	PHE	5.4
30	BH	127	VAL	5.4
31	BI	57	THR	5.4

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
21	AU	17	THR	5.4
43	DU	59	GLY	5.4
51	D3	31	PRO	5.4
51	B3	45	LYS	5.4
23	BA	2107	C	5.4
9	AI	84	ALA	5.4
37	DO	43	GLU	5.4
3	AC	207	VAL	5.4
19	AS	10	PHE	5.3
7	CG	83	ALA	5.3
19	CS	69	HIS	5.3
13	CM	61	GLU	5.3
51	B3	43	CYS	5.3
10	CJ	39	PRO	5.3
34	BL	150	ALA	5.3
23	DA	362	U	5.3
2	CB	71	VAL	5.2
6	AF	91	VAL	5.2
13	CM	40	ASN	5.2
14	AN	8	GLU	5.2
23	DA	2799	A	5.2
30	BH	128	LEU	5.2
8	CH	131	GLY	5.2
1	AA	80	G	5.2
10	AJ	5	ARG	5.1
9	CI	126	SER	5.1
30	BH	93	THR	5.1
2	AB	187	LEU	5.1
19	AS	9	VAL	5.1
1	AA	1001	G	5.1
51	B3	11	LEU	5.1
2	CB	138	LEU	5.1
30	BH	140	LEU	5.1
23	BA	2104	G	5.1
51	D3	17	LYS	5.1
13	CM	16	ASP	5.1
37	BO	88	ASP	5.1
2	CB	140	HIS	5.0
2	AB	163	PHE	5.0
14	AN	12	ARG	5.0
2	AB	167	PRO	5.0
16	CP	34	GLU	5.0

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Mol	Chain	Res	Type	RSRZ
21	AU	18	TYR	5.0
51	B3	37	ARG	5.0
29	BG	104	GLU	5.0
23	BA	2798	C	5.0
23	BA	1045	A	5.0
29	BG	96	ALA	5.0
37	BO	53	SER	5.0
2	AB	118	LEU	5.0
3	CC	159	GLY	5.0
9	AI	18	PHE	5.0
37	BO	37	ALA	4.9
20	CT	9	ASN	4.9
23	BA	2797	U	4.9
13	AM	117	VAL	4.9
1	CA	1356	G	4.9
9	AI	61	ALA	4.9
10	AJ	34	VAL	4.9
29	BG	58	GLU	4.9
23	BA	2804	C	4.9
19	CS	38	SER	4.9
13	CM	29	ARG	4.9
7	CG	62	PHE	4.9
2	AB	7	VAL	4.9
10	CJ	8	LEU	4.9
30	BH	81	VAL	4.9
2	AB	19	HIS	4.8
12	AL	126	GLU	4.8
10	CJ	33	GLN	4.8
51	B3	31	PRO	4.8
30	BH	1	MET	4.8
20	CT	48	LYS	4.8
44	BV	96	VAL	4.8
44	DV	178	GLU	4.8
31	BI	13	LEU	4.8
44	BV	97	GLU	4.8
13	AM	116	THR	4.8
19	AS	51	VAL	4.8
13	CM	96	LEU	4.8
23	DA	508	G	4.8
10	CJ	6	ILE	4.8
29	BG	82	GLY	4.8
2	CB	144	ARG	4.8

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
14	AN	14	PRO	4.8
23	BA	2802	G	4.7
51	B3	35	GLU	4.7
37	BO	49	VAL	4.7
31	DI	59	ILE	4.7
2	CB	152	PHE	4.7
10	AJ	74	ILE	4.7
3	CC	179	ARG	4.7
30	DH	87	LYS	4.7
3	AC	201	TYR	4.7
44	BV	182	LYS	4.7
7	CG	50	ILE	4.7
6	AF	7	ASN	4.7
7	CG	58	PRO	4.7
23	DA	2211	G	4.7
30	BH	109	ILE	4.7
8	CH	42	GLU	4.7
9	AI	63	ILE	4.6
23	BA	2894	G	4.6
19	AS	69	HIS	4.6
20	CT	64	ASP	4.6
9	AI	3	GLN	4.6
23	BA	2793	G	4.6
1	CA	1112	C	4.6
16	CP	21	VAL	4.6
1	CA	1149	C	4.6
51	B3	23	THR	4.6
43	BU	2	ARG	4.6
14	CN	5	ALA	4.6
23	DA	2804	C	4.6
30	DH	70	GLU	4.6
3	CC	101	LEU	4.5
30	DH	108	THR	4.5
30	BH	91	SER	4.5
37	BO	51	ALA	4.5
20	CT	56	MET	4.5
1	AA	1030	C	4.5
30	DH	105	HIS	4.5
19	AS	40	ILE	4.5
29	BG	94	TYR	4.5
10	AJ	26	ALA	4.5
37	BO	39	ILE	4.5

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Mol	Chain	Res	Type	RSRZ
7	CG	79	ARG	4.5
3	AC	206	GLU	4.5
2	CB	134	GLU	4.4
2	AB	47	THR	4.4
30	DH	130	TYR	4.4
19	CS	70	LYS	4.4
9	AI	4	TYR	4.4
14	CN	14	PRO	4.4
30	DH	106	GLY	4.4
1	CA	87	A	4.4
2	AB	40	HIS	4.4
35	BM	139	GLU	4.4
7	CG	52	GLU	4.4
19	AS	35	SER	4.4
19	CS	81	ARG	4.4
2	AB	43	ASP	4.4
2	AB	101	MET	4.4
2	CB	72	GLY	4.4
19	CS	49	ILE	4.4
13	CM	60	VAL	4.4
12	CL	28	GLY	4.4
30	BH	87	LYS	4.4
14	CN	32	SER	4.4
23	BA	2106	G	4.4
2	CB	141	GLU	4.4
31	BI	56	ASN	4.4
44	BV	127	LYS	4.4
2	CB	99	GLY	4.4
9	CI	105	ASP	4.4
23	DA	2108	C	4.3
14	AN	33	VAL	4.3
30	DH	66	GLU	4.3
9	AI	81	ILE	4.3
22	CV	6188	C	4.3
13	CM	7	VAL	4.3
51	B3	29	ASN	4.3
13	CM	6	GLY	4.3
51	D3	42	TRP	4.3
19	AS	38	SER	4.3
34	DL	121	LYS	4.3
30	BH	145	VAL	4.3
19	CS	50	ALA	4.3

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
51	D3	25	LYS	4.3
16	CP	13	HIS	4.3
19	AS	5	LEU	4.2
37	BO	84	GLN	4.2
9	AI	90	PRO	4.2
31	DI	21	GLN	4.2
1	CA	1036	G	4.2
10	AJ	8	LEU	4.2
19	AS	41	VAL	4.2
2	AB	28	PHE	4.2
16	CP	35	LYS	4.2
1	CA	1002	G	4.2
2	CB	93	VAL	4.2
23	DA	277	C	4.2
2	CB	164	VAL	4.2
2	AB	71	VAL	4.2
10	CJ	41	PRO	4.2
14	AN	55	GLY	4.2
7	AG	104	LEU	4.2
2	AB	202	PRO	4.2
7	CG	84	ASN	4.2
19	CS	29	ARG	4.2
30	DH	140	LEU	4.2
2	CB	167	PRO	4.1
7	AG	85	TYR	4.1
19	AS	60	VAL	4.1
10	CJ	4	ILE	4.1
9	CI	12	GLU	4.1
10	AJ	71	LEU	4.1
14	CN	8	GLU	4.1
43	BU	51	VAL	4.1
1	AA	1492	A	4.1
1	CA	1280	A	4.1
10	AJ	33	GLN	4.1
37	DO	42	ASP	4.1
23	DA	271(D)	U	4.1
29	DG	169	VAL	4.1
3	CC	205	GLY	4.1
18	AR	88	LYS	4.1
14	CN	30	ALA	4.1
44	BV	181	GLU	4.1
7	CG	5	ARG	4.1

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Mol	Chain	Res	Type	RSRZ
29	BG	32	GLU	4.1
10	AJ	68	HIS	4.1
1	CA	90	C	4.1
9	AI	21	PRO	4.1
2	CB	95	GLN	4.1
44	DV	153	SER	4.1
2	CB	27	LYS	4.1
18	AR	31	LEU	4.1
37	DO	19	LYS	4.1
3	AC	152	ILE	4.1
9	AI	65	VAL	4.1
37	BO	106	ARG	4.1
47	DY	16	LEU	4.1
1	CA	81	G	4.1
9	CI	13	ALA	4.1
10	CJ	61	GLU	4.1
13	CM	112	GLY	4.1
9	AI	20	ARG	4.1
31	DI	67	GLY	4.1
37	BO	15	ARG	4.1
31	DI	14	LYS	4.1
7	CG	71	PRO	4.0
28	DF	137	GLU	4.0
51	B3	19	ARG	4.0
2	CB	133	LYS	4.0
3	AC	155	GLY	4.0
21	CU	4	GLY	4.0
5	AE	29	GLY	4.0
29	BG	56	SER	4.0
13	CM	116	THR	4.0
29	BG	124	GLU	4.0
49	B1	56	GLU	4.0
30	DH	91	SER	4.0
2	AB	203	GLY	4.0
8	AH	130	GLY	4.0
13	CM	64	TRP	4.0
23	BA	2105	C	4.0
34	BL	110	TYR	4.0
13	AM	98	VAL	4.0
21	AU	11	GLY	4.0
16	CP	36	ILE	4.0
2	CB	69	LEU	4.0

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
23	DA	2803	C	4.0
9	CI	93	ARG	4.0
30	DH	109	ILE	4.0
51	B3	30	THR	4.0
14	CN	21	TYR	4.0
2	AB	164	VAL	4.0
14	CN	2	ALA	4.0
37	DO	45	GLY	3.9
23	DA	363(A)	G	3.9
14	CN	10	ALA	3.9
10	AJ	100	THR	3.9
1	AA	1138	G	3.9
10	CJ	40	LEU	3.9
1	CA	1248	A	3.9
2	CB	98	LEU	3.9
9	AI	7	THR	3.9
14	CN	7	ILE	3.9
9	AI	101	PHE	3.9
11	AK	117	ASN	3.9
6	AF	90	VAL	3.9
3	CC	19	GLU	3.9
10	AJ	62	HIS	3.9
9	CI	33	PHE	3.9
37	BO	92	TYR	3.9
9	CI	15	ALA	3.9
22	AV	6151	C	3.9
23	DA	2798	C	3.9
2	CB	132	LYS	3.9
23	BA	888	C	3.9
51	D3	10	LEU	3.9
13	AM	101	GLN	3.9
2	AB	201	ILE	3.9
2	AB	227	GLY	3.9
28	BF	2	PRO	3.9
30	DH	110	ASP	3.9
2	AB	12	GLU	3.9
47	DY	12	GLU	3.9
21	CU	8	THR	3.9
27	BE	207	GLY	3.9
13	AM	93	ARG	3.8
6	AF	8	ILE	3.8
1	AA	1224	G	3.8

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Mol	Chain	Res	Type	RSRZ
12	CL	127	ALA	3.8
23	BA	101	G	3.8
8	AH	129	VAL	3.8
21	CU	25	LYS	3.8
29	BG	105	LEU	3.8
30	BH	122	GLU	3.8
23	DA	2797	U	3.8
7	CG	47	CYS	3.8
10	AJ	10	GLY	3.8
4	CD	15	GLU	3.8
23	BA	2187	G	3.8
10	CJ	10	GLY	3.8
30	DH	86	THR	3.8
10	CJ	5	ARG	3.8
19	AS	71	LEU	3.8
37	BO	35	ILE	3.8
2	CB	29	ALA	3.8
10	AJ	59	SER	3.8
19	AS	50	ALA	3.8
37	DO	87	PHE	3.8
23	BA	2805	G	3.8
10	AJ	7	LYS	3.8
29	BG	159	GLU	3.8
30	DH	119	PRO	3.8
13	AM	30	ALA	3.8
23	BA	280	C	3.8
3	AC	148	GLY	3.8
27	DE	207	GLY	3.8
10	CJ	36	GLY	3.8
13	CM	32	GLU	3.8
15	CO	49	ASP	3.8
30	DH	67	ARG	3.8
30	DH	78	THR	3.8
10	CJ	64	GLU	3.7
31	DI	6	ASN	3.7
44	BV	179	ASP	3.7
1	CA	1125	U	3.7
45	BW	76	GLY	3.7
9	CI	128	ARG	3.7
4	CD	182	LYS	3.7
1	CA	664	G	3.7
6	AF	61	LEU	3.7

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Mol	Chain	Res	Type	RSRZ
7	AG	78	ARG	3.7
2	AB	215	LEU	3.7
30	DH	63	ALA	3.7
2	CB	94	ASN	3.7
3	AC	177	THR	3.7
9	AI	127	LYS	3.7
21	AU	21	TYR	3.7
14	CN	9	LYS	3.7
12	AL	127	ALA	3.7
7	CG	49	ILE	3.7
19	CS	61	TYR	3.7
19	CS	47	HIS	3.7
1	CA	193	C	3.7
23	BA	2186	G	3.7
20	CT	84	LEU	3.7
2	AB	31	TYR	3.7
19	AS	33	THR	3.7
20	CT	58	LYS	3.7
2	CB	166	ASP	3.7
23	BA	896	A	3.7
2	CB	218	ALA	3.7
14	AN	10	ALA	3.7
23	DA	2802	G	3.6
14	AN	32	SER	3.6
51	B3	52	VAL	3.6
11	AK	129	SER	3.6
51	B3	27	LYS	3.6
23	DA	2896	C	3.6
1	CA	80	G	3.6
1	CA	1017	G	3.6
8	AH	95	VAL	3.6
2	CB	187	LEU	3.6
13	AM	34	LEU	3.6
47	BY	4	SER	3.6
8	CH	129	VAL	3.6
19	AS	39	THR	3.6
26	BD	204	ALA	3.6
22	CV	6150	U	3.6
30	DH	118	LYS	3.6
44	BV	92	SER	3.6
3	AC	151	VAL	3.6
5	CE	19	MET	3.6

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
31	DI	18	GLU	3.6
3	AC	15	THR	3.6
29	BG	95	ARG	3.6
31	BI	10	LEU	3.6
1	AA	728	A	3.6
3	CC	8	ILE	3.6
16	CP	9	PHE	3.6
2	AB	170	GLU	3.6
30	BH	144	VAL	3.6
10	AJ	67	THR	3.6
2	CB	70	PHE	3.6
22	AV	6153	C	3.6
2	CB	130	ARG	3.6
2	AB	42	ILE	3.6
2	CB	165	VAL	3.6
38	DP	2	ASN	3.6
7	CG	4	ARG	3.6
9	AI	92	TYR	3.6
13	CM	97	PRO	3.6
31	BI	58	LEU	3.6
29	BG	41	MET	3.6
37	BO	86	ALA	3.6
6	AF	6	VAL	3.5
10	AJ	28	ARG	3.5
10	CJ	15	THR	3.5
30	DH	71	ILE	3.5
31	BI	21	GLN	3.5
7	AG	15	ASP	3.5
35	DM	139	GLU	3.5
1	CA	136(B)	C	3.5
2	AB	194	PRO	3.5
30	BH	83	ALA	3.5
34	DL	110	TYR	3.5
9	CI	96	LEU	3.5
51	B3	9	LEU	3.5
23	BA	1044	G	3.5
13	CM	98	VAL	3.5
30	BH	65	ALA	3.5
1	AA	1042	G	3.5
31	DI	22	GLY	3.5
11	AK	128	ALA	3.5
9	AI	19	LEU	3.5

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Mol	Chain	Res	Type	RSRZ
38	BP	25	GLY	3.5
18	AR	34	TYR	3.5
12	AL	49	SER	3.5
28	BF	85	GLY	3.5
1	CA	186(A)	C	3.5
20	CT	69	GLY	3.5
37	BO	68	GLN	3.5
9	CI	124	GLN	3.5
28	BF	171	ALA	3.5
38	DP	1	MET	3.5
51	B3	36	LEU	3.5
13	CM	69	GLU	3.5
10	AJ	88	LEU	3.5
18	AR	47	THR	3.5
30	DH	98	ALA	3.5
30	DH	114	LEU	3.4
6	AF	88	VAL	3.4
10	AJ	70	ARG	3.4
44	BV	189	ALA	3.4
48	BZ	34	GLU	3.4
9	CI	125	TYR	3.4
18	AR	32	ARG	3.4
48	BZ	2	PRO	3.4
11	CK	116	HIS	3.4
31	BI	18	GLU	3.4
13	AM	99	ARG	3.4
51	D3	29	ASN	3.4
6	AF	94	GLN	3.4
28	BF	13	GLU	3.4
2	AB	48	MET	3.4
49	B1	60	GLU	3.4
11	AK	11	LYS	3.4
1	CA	1039	C	3.4
23	BA	2803	C	3.4
20	CT	83	ARG	3.4
46	DX	85	LEU	3.4
2	AB	166	ASP	3.4
19	CS	35	SER	3.4
10	AJ	61	GLU	3.4
3	CC	204	LEU	3.4
30	DH	121	LYS	3.4
41	DS	7	ALA	3.4

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
2	CB	54	THR	3.4
13	AM	96	LEU	3.4
37	DO	23	ARG	3.4
20	CT	63	ILE	3.4
48	BZ	27	GLY	3.4
2	CB	169	LYS	3.4
9	AI	26	VAL	3.4
21	AU	8	THR	3.4
43	BU	52	SER	3.4
3	CC	160	ALA	3.4
7	CG	134	ALA	3.4
8	AH	131	GLY	3.4
10	AJ	6	ILE	3.4
21	CU	7	ARG	3.4
29	DG	170	ARG	3.4
44	BV	91	LEU	3.3
11	AK	109	VAL	3.3
1	CA	1129	C	3.3
6	AF	63	TYR	3.3
23	BA	267	C	3.3
23	BA	279	C	3.3
28	DF	2	PRO	3.3
2	CB	137	ARG	3.3
2	AB	50	GLU	3.3
14	CN	6	LEU	3.3
6	CF	1	MET	3.3
3	AC	11	ARG	3.3
9	AI	80	GLY	3.3
22	CV	6151	C	3.3
29	BG	93	GLY	3.3
23	BA	271(D)	U	3.3
9	AI	79	LEU	3.3
23	DA	899	A	3.3
48	DZ	57	GLU	3.3
4	CD	42	GLN	3.3
13	AM	39	ILE	3.3
13	CM	33	ALA	3.3
37	BO	102	ALA	3.3
7	AG	101	LEU	3.3
7	CG	85	TYR	3.3
7	CG	69	VAL	3.3
9	CI	14	VAL	3.3

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Mol	Chain	Res	Type	RSRZ
23	BA	2190	G	3.3
2	AB	77	ALA	3.3
2	AB	99	GLY	3.3
13	AM	100	GLY	3.3
34	BL	88	LEU	3.3
13	CM	4	ILE	3.3
2	CB	122	PHE	3.3
1	AA	1036	G	3.3
12	AL	48	ASN	3.3
20	CT	65	LYS	3.3
10	CJ	54	PHE	3.3
37	DO	28	VAL	3.3
43	DU	17	SER	3.3
30	BH	129	THR	3.3
44	DV	179	ASP	3.3
1	CA	186(B)	C	3.3
1	CA	1000	A	3.3
7	AG	16	LEU	3.3
28	BF	178	PHE	3.3
30	BH	36	ALA	3.3
37	DO	27	SER	3.3
6	AF	92	LYS	3.3
23	BA	2189	U	3.3
23	DA	2897	U	3.3
19	CS	66	MET	3.3
28	DF	54	GLU	3.3
1	CA	1260	C	3.3
14	AN	19	ARG	3.3
29	BG	43	VAL	3.3
29	BG	90	LYS	3.3
37	BO	36	TYR	3.3
51	D3	30	THR	3.3
21	CU	2	GLY	3.3
10	AJ	3	LYS	3.3
14	AN	56	VAL	3.3
1	AA	723	U	3.2
7	CG	80	VAL	3.2
1	CA	990	C	3.2
4	CD	169	LYS	3.2
3	AC	149	ALA	3.2
35	DM	105	GLU	3.2
37	BO	107	GLU	3.2

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
10	CJ	74	ILE	3.2
3	AC	12	LEU	3.2
2	AB	140	HIS	3.2
10	CJ	65	LEU	3.2
3	AC	166	GLU	3.2
5	CE	80	ILE	3.2
17	CQ	9	VAL	3.2
2	CB	168	THR	3.2
14	CN	23	ARG	3.2
23	BA	359	A	3.2
13	CM	66	LEU	3.2
37	BO	24	LEU	3.2
2	AB	41	ILE	3.2
2	AB	200	ILE	3.2
13	AM	112	GLY	3.2
31	BI	19	ARG	3.2
2	CB	163	PHE	3.2
7	AG	12	LEU	3.2
37	BO	27	SER	3.2
4	CD	62	GLN	3.2
1	AA	1027	C	3.2
1	CA	743	U	3.2
28	BF	76	SER	3.2
16	CP	18	ARG	3.2
9	AI	29	ASN	3.2
2	AB	18	GLY	3.2
4	AD	152	SER	3.2
9	AI	87	GLN	3.2
9	CI	9	ARG	3.2
2	AB	70	PHE	3.2
7	CG	101	LEU	3.2
13	CM	19	LEU	3.2
44	BV	188	ALA	3.2
44	BV	187	ALA	3.2
23	BA	277	C	3.2
2	AB	214	ILE	3.2
2	CB	200	ILE	3.2
12	CL	72	GLU	3.2
30	BH	77	LEU	3.2
3	CC	103	VAL	3.2
14	CN	56	VAL	3.2
1	AA	1031	G	3.2

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Mol	Chain	Res	Type	RSRZ
9	AI	34	ASN	3.2
22	AV	6152	G	3.2
11	CK	128	ALA	3.2
18	AR	76	LEU	3.2
28	BF	26	GLN	3.2
13	AM	29	ARG	3.2
16	CP	8	ARG	3.2
37	DO	41	ASP	3.1
38	DP	136	GLN	3.1
2	CB	12	GLU	3.1
43	DU	2	ARG	3.1
47	DY	14	ARG	3.1
13	CM	34	LEU	3.1
18	AR	78	LEU	3.1
19	AS	70	LYS	3.1
20	CT	57	ARG	3.1
28	BF	21	ARG	3.1
1	AA	1260	C	3.1
1	CA	666	G	3.1
3	CC	156	ARG	3.1
7	CG	22	LEU	3.1
28	BF	172	LEU	3.1
19	AS	12	ASP	3.1
1	CA	91	C	3.1
16	CP	29	ASP	3.1
20	CT	50	GLU	3.1
23	DA	2602	A	3.1
7	AG	5	ARG	3.1
10	CJ	59	SER	3.1
29	DG	101	ARG	3.1
38	BP	2	ASN	3.1
28	DF	69	ALA	3.1
49	B1	63	SER	3.1
14	AN	20	ALA	3.1
34	BL	148	LEU	3.1
2	AB	211	ILE	3.1
29	DG	116	GLU	3.1
30	BH	112	LYS	3.1
1	CA	508	C	3.1
22	CV	6155	C	3.1
47	BY	3	LEU	3.1
1	CA	663	A	3.1

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
30	DH	101	LEU	3.1
3	AC	17	ASP	3.1
44	BV	95	PRO	3.1
1	CA	1533	C	3.1
8	CH	133	LEU	3.1
20	CT	70	SER	3.1
30	DH	5	LEU	3.1
9	CI	101	PHE	3.1
51	D3	28	ARG	3.1
28	DF	146	TYR	3.1
2	AB	218	ALA	3.1
9	CI	18	PHE	3.1
7	CG	66	VAL	3.1
37	DO	88	ASP	3.1
44	BV	178	GLU	3.1
3	AC	10	PHE	3.1
23	BA	1128	A	3.1
14	AN	30	ALA	3.1
1	CA	89	U	3.0
30	DH	116	LEU	3.0
3	AC	2	GLY	3.0
13	AM	43	THR	3.0
19	CS	59	PRO	3.0
28	DF	66	GLN	3.0
18	CR	32	ARG	3.0
9	CI	81	ILE	3.0
21	AU	12	LYS	3.0
2	CB	31	TYR	3.0
7	CG	130	GLY	3.0
34	DL	119	GLU	3.0
10	AJ	87	THR	3.0
34	DL	137	LYS	3.0
43	DU	54	LYS	3.0
8	CH	44	PHE	3.0
9	CI	102	LEU	3.0
49	B1	53	THR	3.0
29	BG	123	PHE	3.0
31	BI	20	ALA	3.0
48	DZ	29	ARG	3.0
37	DO	36	TYR	3.0
8	CH	121	ASP	3.0
12	CL	113	LYS	3.0

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Mol	Chain	Res	Type	RSRZ
13	AM	102	ARG	3.0
4	CD	16	GLY	3.0
8	CH	119	LEU	3.0
19	AS	37	ARG	3.0
8	CH	109	ILE	3.0
34	DL	149	GLU	3.0
30	BH	20	ASP	3.0
20	CT	54	LYS	3.0
3	CC	16	ARG	3.0
23	BA	6	A	3.0
10	AJ	12	ASP	3.0
9	AI	33	PHE	3.0
4	CD	108	LEU	3.0
34	BL	149	GLU	3.0
22	CV	6167	G	3.0
29	DG	59	ARG	3.0
13	CM	41	PRO	3.0
15	CO	2	PRO	3.0
2	AB	96	ARG	3.0
2	AB	111	ARG	3.0
29	BG	161	GLY	3.0
13	CM	9	ILE	3.0
37	BO	42	ASP	3.0
38	DP	91	ARG	3.0
49	B1	59	VAL	3.0
15	CO	51	HIS	3.0
8	AH	102	ARG	3.0
17	CQ	22	LEU	3.0
19	CS	37	ARG	3.0
23	BA	508	G	3.0
23	BA	1107	G	3.0
51	B3	18	ARG	3.0
9	CI	43	ALA	3.0
1	CA	1137	C	3.0
1	CA	1289	A	3.0
28	DF	147	ASP	3.0
34	DL	81	GLN	2.9
13	CM	87	TYR	2.9
48	DZ	8	LEU	2.9
10	AJ	19	SER	2.9
17	CQ	26	GLN	2.9
10	CJ	100	THR	2.9

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Mol	Chain	Res	Type	RSRZ
23	BA	362	U	2.9
29	BG	34	GLU	2.9
43	BU	50	ARG	2.9
18	AR	48	GLY	2.9
37	BO	82	ILE	2.9
6	CF	95	GLU	2.9
22	AV	6177	G	2.9
35	BM	91	GLU	2.9
44	BV	185	GLU	2.9
18	CR	31	LEU	2.9
20	CT	99	LEU	2.9
8	CH	94	TYR	2.9
13	CM	100	GLY	2.9
38	DP	90	GLN	2.9
38	DP	22	PHE	2.9
31	DI	10	LEU	2.9
9	AI	8	GLY	2.9
1	CA	1148	U	2.9
9	CI	63	ILE	2.9
1	CA	972	C	2.9
13	AM	94	ARG	2.9
20	CT	80	ARG	2.9
37	BO	38	GLN	2.9
13	AM	106	ASN	2.9
2	AB	197	VAL	2.9
19	CS	73	GLU	2.9
2	AB	188	ALA	2.9
2	CB	203	GLY	2.9
9	CI	62	TYR	2.9
37	BO	90	GLY	2.9
1	AA	1267	C	2.9
23	BA	2896	C	2.9
7	CG	53	LYS	2.9
3	CC	158	GLY	2.9
13	CM	80	ARG	2.9
4	AD	134	ASP	2.9
37	BO	57	LYS	2.9
48	DZ	10	LYS	2.9
44	BV	163	LEU	2.9
13	AM	113	PRO	2.9
38	BP	1	MET	2.9
6	AF	50	TYR	2.9

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
10	CJ	3	LYS	2.9
45	DW	76	GLY	2.9
30	DH	11	ASN	2.9
37	BO	28	VAL	2.9
44	BV	190	GLU	2.9
49	B1	58	TYR	2.9
9	CI	104	ARG	2.9
10	AJ	38	ILE	2.9
37	DO	40	ILE	2.9
2	AB	16	HIS	2.9
29	BG	111	HIS	2.9
37	BO	25	ARG	2.9
48	BZ	58	VAL	2.9
1	AA	727	G	2.9
1	CA	44	G	2.9
1	CA	727	G	2.9
37	BO	19	LYS	2.9
31	BI	22	GLY	2.9
20	CT	88	VAL	2.9
41	BS	101	SER	2.9
29	BG	40	GLU	2.9
50	B2	48	GLU	2.9
9	CI	85	LEU	2.8
2	AB	160	ASP	2.8
6	AF	9	VAL	2.8
15	CO	46	HIS	2.8
3	AC	197	GLY	2.8
14	CN	51	GLY	2.8
2	AB	44	LEU	2.8
2	CB	118	LEU	2.8
10	CJ	37	PRO	2.8
29	BG	17	VAL	2.8
51	B3	10	LEU	2.8
1	CA	1355	G	2.8
44	DV	88	PHE	2.8
9	AI	15	ALA	2.8
44	DV	27	VAL	2.8
3	CC	66	VAL	2.8
9	CI	97	LYS	2.8
25	BC	35	LYS	2.8
37	BO	65	VAL	2.8
23	BA	1535	U	2.8

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Mol	Chain	Res	Type	RSRZ
10	AJ	4	ILE	2.8
20	AT	65	LYS	2.8
27	DE	194	MET	2.8
8	AH	94	TYR	2.8
29	BG	131	VAL	2.8
7	AG	108	ALA	2.8
1	CA	726	C	2.8
43	DU	89	PHE	2.8
20	CT	87	LYS	2.8
7	AG	134	ALA	2.8
19	CS	42	PRO	2.8
26	DD	112	GLY	2.8
37	BO	41	ASP	2.8
20	CT	106	ALA	2.8
28	DF	76	SER	2.8
7	CG	48	LYS	2.8
7	AG	80	VAL	2.8
16	CP	30	GLY	2.8
48	BZ	28	LEU	2.8
22	AV	6193	A	2.8
23	BA	276	A	2.8
23	BA	2794	C	2.8
10	AJ	23	ILE	2.8
2	AB	152	PHE	2.8
12	CL	31	PHE	2.8
2	CB	102	LEU	2.8
37	BO	73	LEU	2.8
1	CA	103(B)	G	2.8
37	BO	22	GLY	2.8
13	CM	31	LYS	2.8
29	BG	61	HIS	2.8
30	DH	97	ILE	2.8
20	CT	62	LEU	2.8
23	DA	2794	C	2.8
30	DH	35	LEU	2.8
9	AI	83	ARG	2.8
13	CM	21	TYR	2.8
32	BJ	142	ARG	2.8
1	CA	1139	G	2.8
9	CI	17	VAL	2.8
48	DZ	28	LEU	2.8
8	CH	111	ILE	2.8

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
13	CM	57	ARG	2.8
1	AA	1092	A	2.8
1	AA	1319	A	2.8
9	AI	46	ALA	2.8
35	DM	91	GLU	2.8
20	CT	71	THR	2.8
30	BH	96	ASP	2.8
7	AG	79	ARG	2.8
30	DH	125	GLU	2.8
34	DL	89	ALA	2.8
30	DH	145	VAL	2.8
9	CI	16	ARG	2.8
2	CB	68	ILE	2.8
10	CJ	53	PRO	2.8
44	BV	183	LEU	2.7
14	CN	19	ARG	2.7
26	BD	186	GLY	2.7
44	DV	118	GLN	2.7
10	AJ	44	VAL	2.7
20	AT	26	ASN	2.7
29	BG	64	LEU	2.7
30	DH	12	LEU	2.7
13	CM	13	LYS	2.7
29	BG	67	LEU	2.7
29	BG	125	VAL	2.7
23	DA	1026	U	2.7
37	BO	40	ILE	2.7
1	AA	1451	A	2.7
1	CA	994	A	2.7
6	CF	94	GLN	2.7
10	AJ	85	LEU	2.7
4	CD	107	ARG	2.7
7	CG	3	ARG	2.7
3	CC	152	ILE	2.7
27	DE	192	LEU	2.7
44	DV	186	GLU	2.7
3	CC	177	THR	2.7
1	AA	1043	C	2.7
1	AA	1323	G	2.7
3	CC	153	VAL	2.7
23	DA	2805	G	2.7
18	AR	72	ARG	2.7

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
44	DV	182	LYS	2.7
12	CL	91	ASP	2.7
3	CC	6	HIS	2.7
44	BV	162	GLU	2.7
10	AJ	35	SER	2.7
19	AS	36	ARG	2.7
10	CJ	76	ASN	2.7
16	AP	59	TRP	2.7
19	CS	76	PRO	2.7
51	D3	45	LYS	2.7
10	CJ	45	ARG	2.7
41	BS	84	ARG	2.7
2	CB	129	GLU	2.7
18	AR	81	PHE	2.7
9	AI	9	ARG	2.7
23	DA	2182	G	2.7
43	BU	5	MET	2.7
10	AJ	58	ASP	2.7
16	CP	11	SER	2.7
17	CQ	7	THR	2.7
26	BD	24	THR	2.7
48	DZ	34	GLU	2.7
2	CB	202	PRO	2.7
10	CJ	9	ARG	2.7
14	CN	18	VAL	2.7
1	CA	1140	C	2.7
22	AV	6176	C	2.7
10	AJ	64	GLU	2.7
1	CA	1001	G	2.7
2	AB	193	ASP	2.7
2	CB	192	SER	2.7
5	CE	24	ARG	2.7
23	BA	2795	G	2.7
28	DF	34	LEU	2.7
10	CJ	7	LYS	2.7
1	CA	1286	A	2.7
2	AB	231	GLU	2.7
29	BG	97	ARG	2.7
34	DL	5	ASP	2.7
10	AJ	37	PRO	2.7
21	AU	22	ARG	2.7
7	AG	81	GLY	2.7

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Mol	Chain	Res	Type	RSRZ
19	CS	39	THR	2.7
37	BO	48	LEU	2.7
1	CA	1243	C	2.7
9	AI	128	ARG	2.7
23	DA	1919	A	2.7
37	BO	97	ARG	2.7
4	CD	8	VAL	2.7
19	AS	15	LEU	2.7
13	AM	97	PRO	2.7
47	DY	9	GLN	2.7
9	AI	48	GLU	2.6
19	CS	4	SER	2.6
11	AK	118	GLY	2.6
48	DZ	33	GLN	2.6
27	BE	181	LEU	2.6
29	BG	19	VAL	2.6
23	DA	898	C	2.6
6	CF	2	ARG	2.6
30	BH	118	LYS	2.6
2	AB	90	MET	2.6
8	AH	119	LEU	2.6
30	DH	81	VAL	2.6
44	BV	128	VAL	2.6
48	BZ	26	LEU	2.6
38	DP	46	GLU	2.6
38	DP	92	GLY	2.6
44	DV	190	GLU	2.6
2	AB	97	TRP	2.6
1	AA	731	G	2.6
1	AA	791	G	2.6
9	CI	44	VAL	2.6
13	CM	25	ILE	2.6
14	CN	49	HIS	2.6
2	CB	32	ILE	2.6
49	B1	36	VAL	2.6
1	AA	1044	A	2.6
9	AI	5	TYR	2.6
19	CS	44	MET	2.6
30	DH	69	LYS	2.6
1	CA	723	U	2.6
31	DI	20	ALA	2.6
29	BG	112	PRO	2.6

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Mol	Chain	Res	Type	RSRZ
17	CQ	32	TYR	2.6
37	BO	21	THR	2.6
37	BO	50	SER	2.6
7	AG	130	GLY	2.6
10	AJ	98	ILE	2.6
8	AH	113	SER	2.6
16	AP	1	MET	2.6
27	DE	191	ARG	2.6
29	BG	59	ARG	2.6
9	AI	89	ASN	2.6
41	BS	100	THR	2.6
24	DB	41	U	2.6
16	CP	19	ILE	2.6
30	BH	21	VAL	2.6
37	BO	95	HIS	2.6
48	DZ	35	ARG	2.6
13	CM	68	GLY	2.6
14	AN	2	ALA	2.6
3	CC	178	LEU	2.6
9	AI	88	TYR	2.6
14	AN	21	TYR	2.6
47	BY	9	GLN	2.6
1	AA	1371	G	2.6
1	CA	82	U	2.6
20	CT	39	LYS	2.6
23	BA	2808	U	2.6
16	CP	7	ALA	2.6
9	AI	2	GLU	2.6
9	AI	56	LEU	2.6
37	DO	25	ARG	2.6
49	B1	51	TYR	2.6
13	AM	7	VAL	2.5
13	AM	60	VAL	2.5
30	DH	84	GLY	2.6
28	BF	113	ARG	2.5
7	AG	18	TYR	2.5
8	AH	25	ASP	2.5
11	AK	110	ASP	2.5
19	CS	60	VAL	2.5
44	DV	175	VAL	2.5
3	CC	196	LEU	2.5
43	BU	62	GLU	2.5

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Mol	Chain	Res	Type	RSRZ
37	DO	38	GLN	2.5
34	DL	120	ALA	2.5
10	CJ	75	ILE	2.5
36	BN	72	ASP	2.5
21	AU	25	LYS	2.5
20	CT	101	GLY	2.5
23	DA	887	A	2.5
2	AB	46	LYS	2.5
28	DF	152	LEU	2.5
24	DB	8	U	2.5
17	CQ	27	PHE	2.5
2	AB	168	THR	2.5
7	CG	91	VAL	2.5
20	CT	79	ARG	2.5
7	CG	46	ALA	2.5
14	AN	7	ILE	2.5
23	BA	274	G	2.5
7	AG	20	ASP	2.5
12	CL	111	ASP	2.5
2	CB	77	ALA	2.5
2	CB	131	PRO	2.5
15	CO	52	SER	2.5
21	AU	19	GLY	2.5
31	DI	65	GLU	2.5
9	CI	41	VAL	2.5
17	CQ	23	VAL	2.5
43	BU	45	VAL	2.5
9	AI	6	GLY	2.5
12	CL	70	PRO	2.5
28	BF	34	LEU	2.5
7	CG	54	THR	2.5
29	BG	106	THR	2.5
1	AA	792	A	2.5
1	CA	1035	A	2.5
2	AB	226	ARG	2.5
7	AG	44	TYR	2.5
24	DB	53	A	2.5
2	CB	81	VAL	2.5
8	CH	110	ALA	2.5
11	CK	65	ALA	2.5
6	AF	60	PHE	2.5
49	B1	50	THR	2.5

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
3	AC	150	LYS	2.5
1	CA	136(A)	C	2.5
2	AB	54	THR	2.5
26	DD	159	HIS	2.5
28	DF	35	GLU	2.5
2	CB	215	LEU	2.5
13	CM	95	GLY	2.5
18	CR	66	LEU	2.5
7	AG	84	ASN	2.5
12	CL	44	PRO	2.5
7	AG	105	VAL	2.5
1	AA	1026	G	2.5
29	BG	98	LEU	2.5
30	DH	77	LEU	2.5
2	AB	223	ILE	2.5
35	DM	24	GLY	2.5
3	CC	73	PRO	2.5
18	CR	63	GLN	2.5
34	BL	9	ASN	2.5
6	CF	4	TYR	2.5
9	AI	36	TYR	2.5
1	AA	999	U	2.5
3	AC	54	ARG	2.5
20	CT	76	ALA	2.5
48	BZ	5	LYS	2.5
7	CG	61	VAL	2.5
24	DB	54	G	2.5
31	DI	16	ASN	2.5
27	DE	156	LEU	2.5
2	AB	137	ARG	2.5
2	CB	57	PHE	2.5
2	CB	42	ILE	2.5
6	AF	5	GLU	2.5
30	DH	79	ILE	2.5
4	CD	110	PHE	2.5
9	AI	126	SER	2.5
1	CA	1249	C	2.5
30	BH	19	VAL	2.5
49	D1	59	VAL	2.5
19	AS	42	PRO	2.5
30	DH	9	LEU	2.5
4	CD	17	VAL	2.5

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
30	BH	123	LEU	2.4
8	AH	96	GLY	2.4
18	AR	49	LYS	2.4
30	BH	4	ILE	2.4
37	DO	22	GLY	2.4
2	AB	206	ASP	2.4
23	BA	1733	G	2.4
30	BH	94	ALA	2.4
4	CD	24	GLU	2.4
7	AG	113	GLU	2.4
13	CM	73	GLU	2.4
34	DL	82	GLY	2.4
35	DM	65	PHE	2.4
28	DF	149	VAL	2.4
10	AJ	65	LEU	2.4
21	AU	5	ASP	2.4
22	CV	6156	A	2.4
20	AT	71	THR	2.4
1	AA	1116	C	2.4
3	CC	2	GLY	2.4
29	BG	81	GLU	2.4
48	BZ	57	GLU	2.4
3	CC	149	ALA	2.4
29	DG	103	LEU	2.4
43	BU	53	PRO	2.4
18	AR	71	LYS	2.4
44	DV	183	LEU	2.4
1	CA	43	C	2.4
1	CA	88	C	2.4
1	CA	979	C	2.4
23	BA	1920	C	2.4
10	CJ	63	PHE	2.4
1	CA	1124	G	2.4
8	CH	95	VAL	2.4
23	BA	10	G	2.4
23	BA	2893	G	2.4
2	CB	145	LEU	2.4
18	CR	64	ARG	2.4
19	CS	65	ASN	2.4
14	CN	15	LYS	2.4
17	CQ	41	LYS	2.4
23	BA	2188	C	2.4

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
48	BZ	33	GLN	2.4
36	DN	109	ALA	2.4
1	CA	1220	G	2.4
48	BZ	29	ARG	2.4
1	AA	994	A	2.4
2	AB	130	ARG	2.4
11	CK	82	VAL	2.4
20	AT	103	GLY	2.4
23	BA	2895	U	2.4
28	DF	92	VAL	2.4
30	BH	142	VAL	2.4
44	DV	181	GLU	2.4
50	D2	48	GLU	2.4
2	AB	102	LEU	2.4
8	AH	35	ILE	2.4
1	CA	1003	G	2.4
23	BA	1112	G	2.4
2	AB	33	TYR	2.4
5	CE	118	ILE	2.4
2	CB	190	THR	2.4
9	AI	14	VAL	2.4
12	AL	46	LYS	2.4
19	CS	33	THR	2.4
45	BW	42	GLY	2.4
3	CC	193	TYR	2.4
13	AM	4	ILE	2.4
20	CT	38	LYS	2.4
20	CT	61	SER	2.4
23	BA	2792	G	2.4
28	DF	70	VAL	2.4
37	BO	23	ARG	2.4
38	BP	106	SER	2.4
16	AP	7	ALA	2.4
19	CS	78	ARG	2.4
28	BF	19	LEU	2.4
29	BG	68	THR	2.4
1	CA	1418	A	2.4
22	AV	6189	A	2.4
23	DA	1916	A	2.4
48	BZ	35	ARG	2.4
2	AB	204	ASN	2.4
5	CE	81	GLU	2.4

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
18	AR	69	THR	2.4
29	DG	104	GLU	2.4
28	DF	153	ARG	2.4
2	CB	92	TYR	2.4
2	CB	135	GLN	2.4
10	CJ	55	LYS	2.4
2	AB	136	VAL	2.4
5	CE	98	THR	2.4
1	CA	744	C	2.3
6	AF	4	TYR	2.4
19	CS	67	VAL	2.4
53	B5	31	HIS	2.4
23	DA	897	C	2.3
2	AB	76	GLN	2.3
26	DD	186	GLY	2.3
8	CH	113	SER	2.3
10	CJ	95	GLU	2.3
17	CQ	36	ILE	2.3
49	D1	57	ILE	2.3
19	AS	34	TRP	2.3
1	AA	196	A	2.3
2	AB	100	GLY	2.3
7	AG	26	PHE	2.3
7	AG	43	PHE	2.3
22	AV	6167	G	2.3
2	AB	81	VAL	2.3
44	BV	27	VAL	2.3
13	CM	63	THR	2.3
1	CA	1150	U	2.3
3	AC	170	GLN	2.3
19	CS	32	LYS	2.3
2	CB	201	ILE	2.3
14	CN	16	PHE	2.3
23	BA	1111	A	2.3
3	AC	147	LYS	2.3
12	AL	110	LYS	2.3
28	BF	173	LEU	2.3
41	BS	82	LEU	2.3
1	CA	102(B)	C	2.3
1	CA	1127	G	2.3
12	CL	18	ARG	2.3
23	DA	2334	G	2.3

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
23	DA	2793	G	2.3
28	DF	115	ARG	2.3
43	DU	86	ARG	2.3
25	BC	36	PRO	2.3
16	CP	17	TYR	2.3
16	CP	32	TYR	2.3
21	AU	20	LYS	2.3
30	BH	5	LEU	2.3
2	AB	72	GLY	2.3
4	CD	47	ARG	2.3
1	AA	1252	A	2.3
1	AA	1318	A	2.3
31	DI	8	GLU	2.3
13	AM	56	LEU	2.3
18	AR	44	LEU	2.3
48	DZ	53	LEU	2.3
18	AR	22	VAL	2.3
19	AS	58	VAL	2.3
23	DA	1917	U	2.3
32	BJ	73	ASP	2.3
34	DL	114	ILE	2.3
7	CG	10	ARG	2.3
10	CJ	60	ARG	2.3
34	BL	138	LEU	2.3
37	DO	37	ALA	2.3
2	CB	214	ILE	2.3
9	AI	37	PHE	2.3
2	AB	51	LEU	2.3
2	AB	159	PRO	2.3
2	CB	142	LEU	2.3
18	CR	79	LEU	2.3
20	CT	68	LYS	2.3
31	DI	13	LEU	2.3
1	CA	1138	G	2.3
10	AJ	63	PHE	2.3
51	D3	27	LYS	2.3
4	CD	10	ARG	2.3
6	AF	69	GLU	2.3
7	CG	78	ARG	2.3
38	DP	36	GLU	2.3
48	DZ	30	ARG	2.3
4	CD	183	GLY	2.3

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Mol	Chain	Res	Type	RSRZ
37	BO	81	GLY	2.3
1	CA	194	C	2.3
23	DA	1537	C	2.3
28	DF	88	ILE	2.3
1	CA	191(G)	G	2.3
1	CA	742	G	2.3
23	BA	11	G	2.3
2	CB	193	ASP	2.3
31	DI	12	THR	2.3
38	DP	26	ASP	2.3
29	BG	115	VAL	2.3
37	BO	85	VAL	2.3
27	DE	155	LEU	2.3
4	CD	4	TYR	2.3
1	CA	1115	C	2.3
2	AB	230	VAL	2.3
10	CJ	23	ILE	2.3
19	AS	32	LYS	2.3
30	BH	37	VAL	2.3
37	DO	11	LYS	2.3
7	AG	37	ASN	2.3
23	BA	226	G	2.3
37	BO	16	ASN	2.3
48	DZ	4	LEU	2.3
2	CB	26	PRO	2.3
2	AB	93	VAL	2.3
13	CM	54	VAL	2.3
28	DF	39	ILE	2.3
34	BL	7	ARG	2.3
2	CB	59	GLU	2.3
3	CC	169	ALA	2.3
19	CS	43	GLU	2.3
7	AG	88	PRO	2.3
18	AR	77	GLY	2.3
21	CU	23	PRO	2.3
34	BL	137	LYS	2.3
37	BO	83	LYS	2.3
29	BG	89	ILE	2.3
1	CA	975	A	2.3
17	CQ	24	GLU	2.3
8	AH	128	GLY	2.3
10	AJ	24	VAL	2.3

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Mol	Chain	Res	Type	RSRZ
23	BA	1919	A	2.3
11	CK	115	PRO	2.3
11	CK	117	ASN	2.3
30	DH	103	ARG	2.3
1	CA	1037	C	2.3
2	AB	213	LEU	2.3
13	CM	18	ALA	2.3
13	CM	23	TYR	2.3
10	CJ	96	ILE	2.3
32	BJ	56	LEU	2.2
1	CA	1285	A	2.2
2	AB	55	PHE	2.2
9	AI	125	TYR	2.2
23	DA	614	U	2.2
29	DG	60	ARG	2.2
46	BX	20	ARG	2.2
16	CP	41	PRO	2.2
2	AB	122	PHE	2.2
1	CA	509	A	2.2
8	AH	112	LEU	2.2
11	CK	110	ASP	2.2
44	BV	28	MET	2.2
12	CL	30	PRO	2.2
13	AM	111	LYS	2.2
3	CC	59	ARG	2.2
7	AG	22	LEU	2.2
1	CA	85	U	2.2
27	DE	193	VAL	2.2
1	AA	1029	G	2.2
2	CB	121	LEU	2.2
3	AC	13	GLY	2.2
20	CT	66	ALA	2.2
23	BA	2181	G	2.2
23	BA	2807	G	2.2
23	DA	2795	G	2.2
44	BV	113	ALA	2.2
20	CT	35	THR	2.2
18	AR	79	LEU	2.2
21	AU	2	GLY	2.2
27	DE	27	GLU	2.2
30	BH	64	GLU	2.2
2	CB	219	VAL	2.2

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Mol	Chain	Res	Type	RSRZ
12	AL	111	ASP	2.2
14	CN	26	ARG	2.2
19	CS	46	GLY	2.2
20	AT	19	SER	2.2
27	BE	144	LYS	2.2
47	BY	12	GLU	2.2
10	AJ	76	ASN	2.2
7	AG	11	GLN	2.2
13	CM	10	PRO	2.2
1	CA	1330	U	2.2
11	AK	17	GLY	2.2
13	CM	115	LYS	2.2
28	BF	120	LEU	2.2
31	DI	9	LEU	2.2
42	BT	90	GLU	2.2
1	AA	934	C	2.2
1	AA	1338	G	2.2
25	BC	97	TYR	2.2
28	BF	168	GLU	2.2
4	CD	104	VAL	2.2
4	AD	161	ASN	2.2
1	CA	1329	A	2.2
1	CA	1363	A	2.2
10	AJ	18	ALA	2.2
20	CT	53	LEU	2.2
29	BG	103	LEU	2.2
44	BV	164	ALA	2.2
47	BY	13	ALA	2.2
10	AJ	89	ASP	2.2
36	DN	107	ASP	2.2
21	CU	12	LYS	2.2
6	AF	10	LEU	2.2
13	CM	62	ASN	2.2
20	CT	72	LEU	2.2
28	BF	25	TYR	2.2
29	BG	29	PRO	2.2
38	BP	64	ARG	2.2
47	BY	1	MET	2.2
29	BG	25	LYS	2.2
5	AE	14	ARG	2.2
20	CT	95	ALA	2.2
1	AA	177	C	2.2

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Mol	Chain	Res	Type	RSRZ
1	CA	1354	C	2.2
14	AN	9	LYS	2.2
19	CS	28	LYS	2.2
20	CT	85	MET	2.2
5	AE	5	ASP	2.2
23	DA	1465	G	2.2
17	CQ	6	LEU	2.2
18	CR	50	ILE	2.2
2	CB	171	ALA	2.2
22	AV	6178	A	2.2
39	BQ	90	VAL	2.2
1	CA	1267	C	2.2
9	CI	92	TYR	2.2
14	CN	29	ARG	2.2
44	BV	87	ASP	2.2
14	CN	11	LYS	2.2
46	BX	38	SER	2.2
7	CG	13	GLN	2.2
20	CT	42	GLN	2.2
23	BA	361	G	2.2
29	BG	169	VAL	2.2
1	AA	733	A	2.2
1	AA	1363	A	2.2
29	DG	102	ALA	2.2
30	DH	83	ALA	2.2
7	AG	71	PRO	2.2
1	AA	1025	U	2.1
11	CK	12	ARG	2.1
14	AN	23	ARG	2.1
19	CS	57	HIS	2.1
24	DB	88	C	2.1
36	BN	11	ASN	2.1
3	CC	9	GLY	2.1
16	CP	15	PRO	2.1
17	CQ	58	GLU	2.1
28	DF	32	PRO	2.1
29	BG	38	SER	2.1
48	DZ	2	PRO	2.1
1	AA	978	A	2.1
1	AA	1493	A	2.1
6	AF	86	ARG	2.1
31	DI	17	LEU	2.1

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
8	CH	93	VAL	2.1
1	CA	732	C	2.1
1	CA	1100	C	2.1
23	DA	2402	C	2.1
30	DH	7	GLU	2.1
1	AA	1002	G	2.1
10	CJ	94	VAL	2.1
10	CJ	48	THR	2.1
23	DA	7	G	2.1
2	AB	121	LEU	2.1
7	AG	8	GLU	2.1
7	AG	59	LEU	2.1
10	AJ	39	PRO	2.1
43	BU	91	GLU	2.1
1	CA	1257	U	2.1
2	AB	36	ARG	2.1
3	CC	202	ILE	2.1
8	CH	24	THR	2.1
37	BO	30	ARG	2.1
9	AI	49	PRO	2.1
39	BQ	91	ASP	2.1
1	AA	103(B)	G	2.1
8	CH	61	VAL	2.1
4	CD	61	LYS	2.1
4	CD	181	MET	2.1
16	CP	22	THR	2.1
20	AT	21	LYS	2.1
12	AL	47	PRO	2.1
23	BA	2103	C	2.1
28	BF	22	ARG	2.1
37	BO	89	ARG	2.1
2	AB	207	ALA	2.1
2	CB	149	LEU	2.1
13	CM	84	ILE	2.1
16	CP	1	MET	2.1
1	AA	977	A	2.1
1	CA	728	A	2.1
1	CA	1026	G	2.1
2	AB	190	THR	2.1
23	BA	360	G	2.1
23	BA	144(B)	A	2.1
1	AA	1351	U	2.1

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Mol	Chain	Res	Type	RSRZ
9	CI	120	ARG	2.1
16	AP	52	ASP	2.1
16	CP	6	LEU	2.1
41	BS	102	HIS	2.1
2	AB	35	GLU	2.1
8	AH	127	LEU	2.1
1	AA	1257	U	2.1
1	CA	1251	A	2.1
23	BA	2378	A	2.1
1	CA	1224	G	2.1
1	CA	63	C	2.1
1	CA	1030	C	2.1
23	DA	2183	C	2.1
51	B3	34	LEU	2.1
3	CC	56	ASP	2.1
20	CT	46	GLU	2.1
20	CT	105	SER	2.1
11	AK	119	CYS	2.1
28	BF	141	PHE	2.1
29	BG	86	GLU	2.1
1	CA	610	G	2.1
1	CA	1370	G	2.1
8	AH	97	VAL	2.1
31	DI	5	ARG	2.1
37	DO	15	ARG	2.1
1	CA	136	C	2.1
2	AB	80	ILE	2.1
10	CJ	16	LEU	2.1
44	BV	160	GLY	2.1
10	AJ	20	ALA	2.1
31	DI	11	ALA	2.1
19	AS	43	GLU	2.1
44	BV	31	ARG	2.1
23	BA	887	A	2.1
23	BA	1785	A	2.1
28	DF	27	ASN	2.1
48	BZ	52	HIS	2.1
13	CM	83	ASP	2.1
1	CA	976	G	2.1
1	CA	1253	G	2.1
49	D1	37	PRO	2.1
1	AA	995	C	2.1

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Mol	Chain	Res	Type	RSRZ
1	CA	501	C	2.1
1	CA	1116	C	2.1
1	CA	968	A	2.1
8	CH	135	CYS	2.1
23	DA	1536	A	2.1
38	BP	24	PRO	2.1
29	BG	129	THR	2.1
2	CB	66	GLY	2.1
18	CR	40	LEU	2.1
20	AT	102	GLY	2.1
1	AA	1317	C	2.1
1	CA	667	G	2.1
35	DM	63	LYS	2.1
18	CR	82	THR	2.0
9	AI	86	VAL	2.0
44	DV	87	ASP	2.0
48	DZ	9	VAL	2.0
1	CA	1209	C	2.0
4	CD	11	LEU	2.0
27	BE	156	LEU	2.0
29	BG	53	GLU	2.0
14	CN	50	LYS	2.0
16	CP	33	ILE	2.0
43	BU	41	GLY	2.0
19	AS	11	VAL	2.0
29	BG	160	LYS	2.0
49	D1	58	TYR	2.0
4	AD	181	MET	2.0
34	BL	87	ASP	2.0
25	BC	83	GLU	2.0
3	CC	184	TYR	2.0
9	CI	100	GLY	2.0
13	AM	95	GLY	2.0
16	CP	37	GLY	2.0
19	AS	72	GLY	2.0
1	AA	1112	C	2.0
17	CQ	28	PRO	2.0
8	CH	112	LEU	2.0
22	AV	6175	G	2.0
48	DZ	54	VAL	2.0
52	B4	46	VAL	2.0
9	CI	82	ALA	2.0

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Mol	Chain	Res	Type	RSRZ
10	AJ	60	ARG	2.0
18	CR	78	LEU	2.0
28	BF	176	LEU	2.0
9	CI	5	TYR	2.0
12	AL	94	GLY	2.0
16	CP	14	ASN	2.0
1	AA	1137	C	2.0
3	CC	195	VAL	2.0
23	BA	974(B)	C	2.0
7	AG	122	HIS	2.0
43	BU	6	HIS	2.0
1	CA	388	G	2.0
20	CT	81	LYS	2.0
2	AB	45	GLN	2.0
29	DG	167	GLU	2.0
17	CQ	57	VAL	2.0
30	BH	120	ILE	2.0
23	DA	363(B)	A	2.0
43	BU	83	THR	2.0
45	BW	75	LEU	2.0
2	CB	159	PRO	2.0
9	CI	49	PRO	2.0
19	AS	31	ILE	2.0
46	BX	30	VAL	2.0
1	CA	668	G	2.0
1	CA	1186	G	2.0
2	AB	88	ALA	2.0
10	AJ	56	HIS	2.0
13	CM	90	LEU	2.0
13	CM	94	ARG	2.0
30	BH	114	LEU	2.0
7	CG	43	PHE	2.0
32	DJ	74	PHE	2.0
38	DP	27	THR	2.0

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

There are no non-standard protein/DNA/RNA residues in this entry.

## 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, 95<sup>th</sup> 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(Å <sup>2</sup> )	Q<0.9
54	MG	DA	3321	1/1	0.91	0.72	63.58	110,110,110,110	0
54	MG	AA	1613	1/1	0.96	0.63	60.65	55,55,55,55	0
54	MG	AA	1820	1/1	0.47	1.17	60.51	94,94,94,94	0
54	MG	DA	3646	1/1	0.87	0.67	57.35	45,45,45,45	0
54	MG	BA	2926	1/1	0.80	0.69	55.26	50,50,50,50	0
54	MG	BA	3451	1/1	0.78	0.94	54.98	47,47,47,47	0
54	MG	DA	3689	1/1	0.80	1.08	47.20	67,67,67,67	0
54	MG	DA	3178	1/1	0.88	0.66	42.10	54,54,54,54	0
54	MG	BA	2957	1/1	0.85	0.84	41.63	75,75,75,75	0
54	MG	BA	3040	1/1	0.75	1.23	38.37	76,76,76,76	0
54	MG	BA	3425	1/1	0.86	0.67	38.34	85,85,85,85	0
54	MG	DA	3072	1/1	0.87	0.72	37.98	80,80,80,80	0
54	MG	AA	1860	1/1	0.73	0.86	37.30	138,138,138,138	0
54	MG	BA	3643	1/1	0.68	0.84	34.13	92,92,92,92	0
54	MG	DA	3020	1/1	0.92	0.73	33.61	47,47,47,47	0
54	MG	BA	3239	1/1	0.90	0.80	32.70	62,62,62,62	0
54	MG	BA	3060	1/1	0.91	1.14	31.24	110,110,110,110	0
54	MG	BA	3003	1/1	0.78	0.47	30.71	55,55,55,55	0
54	MG	AA	1800	1/1	0.73	1.10	30.07	69,69,69,69	0
54	MG	BA	2929	1/1	0.92	0.69	29.83	50,50,50,50	0
54	MG	DA	3667	1/1	0.90	0.76	29.82	59,59,59,59	0
54	MG	BA	3073	1/1	0.88	0.34	28.99	63,63,63,63	0
54	MG	DA	3103	1/1	0.63	0.67	28.48	58,58,58,58	0
54	MG	AA	1760	1/1	0.91	0.77	28.16	88,88,88,88	0
54	MG	AA	1846	1/1	0.79	0.40	27.84	88,88,88,88	0
54	MG	DA	3227	1/1	0.93	0.49	27.80	43,43,43,43	0
54	MG	DA	3063	1/1	0.88	0.69	26.99	126,126,126,126	0
54	MG	DA	3629	1/1	0.96	0.77	25.85	89,89,89,89	0
54	MG	DA	3065	1/1	0.88	0.41	25.80	61,61,61,61	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
54	MG	BA	3340	1/1	0.95	0.58	25.78	85,85,85,85	0
54	MG	DA	3698	1/1	0.72	0.66	25.45	61,61,61,61	0
54	MG	BA	3196	1/1	0.85	0.86	25.32	80,80,80,80	0
54	MG	DA	3603	1/1	0.79	0.69	25.30	64,64,64,64	0
54	MG	BA	2968	1/1	0.89	0.53	25.00	47,47,47,47	0
54	MG	BA	3136	1/1	0.79	0.36	24.64	64,64,64,64	0
54	MG	BA	3048	1/1	0.90	0.69	24.25	70,70,70,70	0
54	MG	DA	2971	1/1	0.94	0.68	23.94	38,38,38,38	0
54	MG	DA	3080	1/1	0.88	0.78	23.25	83,83,83,83	0
54	MG	DA	3650	1/1	0.83	0.43	22.80	55,55,55,55	0
54	MG	DA	2953	1/1	0.81	0.61	22.72	46,46,46,46	0
54	MG	DA	3637	1/1	0.91	0.47	22.35	15,15,15,15	0
54	MG	BA	3023	1/1	0.86	0.70	22.18	57,57,57,57	0
54	MG	CA	1665	1/1	0.77	1.01	22.02	71,71,71,71	0
54	MG	DA	3013	1/1	0.75	0.90	21.61	52,52,52,52	0
54	MG	DA	3472	1/1	0.95	0.66	21.14	95,95,95,95	0
54	MG	BA	3570	1/1	0.66	0.53	20.73	67,67,67,67	0
54	MG	BA	3159	1/1	0.88	0.69	20.12	59,59,59,59	0
54	MG	BA	3070	1/1	0.74	0.39	20.06	85,85,85,85	0
54	MG	AA	1667	1/1	0.91	0.85	20.06	78,78,78,78	0
54	MG	AA	1724	1/1	0.83	0.35	19.96	79,79,79,79	0
54	MG	DA	2957	1/1	0.90	0.63	19.49	40,40,40,40	0
54	MG	DA	3333	1/1	0.75	0.51	19.43	90,90,90,90	0
54	MG	AA	1620	1/1	0.84	0.58	18.63	56,56,56,56	0
54	MG	BA	3020	1/1	0.85	0.60	18.53	88,88,88,88	0
54	MG	DA	2986	1/1	0.87	0.59	18.15	41,41,41,41	0
54	MG	DA	3177	1/1	0.96	0.59	18.11	46,46,46,46	0
54	MG	DA	3633	1/1	0.86	0.43	17.83	83,83,83,83	0
54	MG	BA	3455	1/1	0.99	0.28	17.39	32,32,32,32	0
54	MG	CA	1676	1/1	0.94	0.49	17.34	90,90,90,90	0
54	MG	BA	3626	1/1	0.81	0.60	17.18	64,64,64,64	0
54	MG	BA	3088	1/1	0.95	0.50	16.97	57,57,57,57	0
54	MG	DA	3018	1/1	0.97	0.44	16.78	60,60,60,60	0
54	MG	BA	3021	1/1	0.90	0.53	16.62	66,66,66,66	0
54	MG	DA	3563	1/1	0.92	0.96	16.47	95,95,95,95	0
54	MG	DA	3144	1/1	0.94	0.32	16.36	69,69,69,69	0
54	MG	DA	3684	1/1	0.98	0.37	16.26	69,69,69,69	0
54	MG	BA	2911	1/1	0.92	0.63	16.07	33,33,33,33	0
54	MG	DA	2923	1/1	0.98	0.46	15.97	12,12,12,12	0
54	MG	BA	3058	1/1	0.85	0.48	15.62	78,78,78,78	0
54	MG	BA	3432	1/1	0.97	0.49	15.52	12,12,12,12	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
54	MG	CA	1875	1/1	0.96	0.64	15.47	63,63,63,63	0
54	MG	BA	3008	1/1	0.83	0.36	15.28	70,70,70,70	0
54	MG	DA	2954	1/1	0.97	0.59	15.08	41,41,41,41	0
54	MG	AA	1631	1/1	0.59	0.43	15.05	76,76,76,76	0
54	MG	DA	3193	1/1	0.82	0.32	15.00	86,86,86,86	0
54	MG	BA	2963	1/1	0.84	0.70	14.97	34,34,34,34	0
54	MG	DA	2947	1/1	0.89	0.48	14.69	37,37,37,37	0
54	MG	BA	2906	1/1	0.98	0.50	14.45	24,24,24,24	0
54	MG	AA	1606	1/1	0.94	0.53	14.29	62,62,62,62	0
54	MG	DC	302	1/1	0.93	0.41	13.88	164,164,164,164	0
54	MG	BA	2939	1/1	0.92	0.52	13.76	47,47,47,47	0
54	MG	BA	3556	1/1	0.92	0.55	13.70	77,77,77,77	0
54	MG	CA	1619	1/1	0.95	0.52	13.66	62,62,62,62	0
54	MG	AA	1664	1/1	0.68	0.44	13.60	80,80,80,80	0
54	MG	DA	3863	1/1	0.84	0.94	13.55	104,104,104,104	0
54	MG	BA	3461	1/1	0.92	0.50	13.45	48,48,48,48	0
54	MG	DA	2913	1/1	0.95	0.67	13.19	21,21,21,21	0
54	MG	DA	3057	1/1	0.83	0.36	13.02	79,79,79,79	0
54	MG	BA	2924	1/1	0.96	0.44	13.02	48,48,48,48	0
54	MG	DA	2901	1/1	0.96	0.53	12.70	9,9,9,9	0
54	MG	BA	3072	1/1	0.94	0.33	12.17	57,57,57,57	0
54	MG	BA	2903	1/1	0.95	0.90	12.11	22,22,22,22	0
54	MG	DA	3154	1/1	0.80	0.39	12.05	50,50,50,50	0
54	MG	BA	3227	1/1	0.94	0.67	11.88	61,61,61,61	0
54	MG	DA	3038	1/1	0.87	0.42	11.71	44,44,44,44	0
54	MG	BA	3122	1/1	0.93	0.24	11.70	56,56,56,56	0
54	MG	BA	2913	1/1	0.94	0.60	11.67	27,27,27,27	0
54	MG	DA	2907	1/1	0.95	0.57	11.53	17,17,17,17	0
54	MG	BA	2920	1/1	0.93	0.41	11.45	24,24,24,24	0
54	MG	DA	2928	1/1	0.93	0.49	11.42	31,31,31,31	0
54	MG	DA	2985	1/1	0.85	0.36	11.42	40,40,40,40	0
54	MG	DA	3511	1/1	0.83	0.51	11.38	52,52,52,52	0
54	MG	AA	1711	1/1	0.91	0.34	11.34	58,58,58,58	0
54	MG	DA	3534	1/1	0.89	0.32	11.31	83,83,83,83	0
54	MG	BA	2954	1/1	0.88	0.27	11.14	73,73,73,73	0
54	MG	AA	1651	1/1	0.93	0.34	10.74	56,56,56,56	0
54	MG	DB	207	1/1	0.97	0.29	10.69	58,58,58,58	0
54	MG	AA	1626	1/1	0.81	0.23	10.68	88,88,88,88	0
54	MG	BA	2916	1/1	0.95	0.47	10.67	37,37,37,37	0
54	MG	DA	3742	1/1	0.94	0.41	10.65	47,47,47,47	0
54	MG	DA	3655	1/1	0.77	0.35	10.59	60,60,60,60	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
54	MG	DA	3733	1/1	0.63	0.49	10.58	84,84,84,84	0
54	MG	DA	3226	1/1	0.85	0.41	10.35	86,86,86,86	0
54	MG	BA	3597	1/1	0.88	0.28	10.29	56,56,56,56	0
54	MG	BA	2934	1/1	0.94	0.53	10.15	33,33,33,33	0
54	MG	DA	3160	1/1	0.96	0.41	10.09	53,53,53,53	0
54	MG	DA	2902	1/1	0.95	0.55	10.06	17,17,17,17	0
54	MG	AA	1647	1/1	0.96	0.44	9.98	63,63,63,63	0
54	MG	DA	2918	1/1	0.98	0.56	9.89	25,25,25,25	0
54	MG	CA	1680	1/1	0.83	0.41	9.83	63,63,63,63	0
54	MG	DA	2982	1/1	0.94	0.26	9.81	39,39,39,39	0
54	MG	DA	3569	1/1	0.93	0.48	9.71	36,36,36,36	0
54	MG	BA	2907	1/1	0.99	0.35	9.64	15,15,15,15	0
54	MG	BA	3049	1/1	0.91	0.54	9.56	58,58,58,58	0
54	MG	BA	2902	1/1	0.97	0.51	9.46	12,12,12,12	0
54	MG	AA	1628	1/1	0.91	0.45	9.42	86,86,86,86	0
54	MG	CA	1602	1/1	0.86	0.51	9.41	43,43,43,43	0
54	MG	DA	3161	1/1	0.95	0.45	9.39	70,70,70,70	0
54	MG	AA	1685	1/1	0.82	0.39	9.28	64,64,64,64	0
54	MG	BA	2922	1/1	0.97	0.53	9.28	38,38,38,38	0
54	MG	CA	1606	1/1	0.98	0.46	9.26	57,57,57,57	0
54	MG	DA	2908	1/1	0.96	0.46	9.15	20,20,20,20	0
54	MG	BA	3172	1/1	0.89	0.26	9.13	74,74,74,74	0
54	MG	CA	1601	1/1	0.92	0.40	9.02	60,60,60,60	0
54	MG	DA	3024	1/1	0.71	0.37	8.96	58,58,58,58	0
54	MG	DA	2925	1/1	0.97	0.40	8.86	34,34,34,34	0
54	MG	DA	3035	1/1	0.95	0.61	8.86	43,43,43,43	0
54	MG	AA	1681	1/1	0.92	0.41	8.70	65,65,65,65	0
54	MG	DA	3859	1/1	0.94	0.40	8.48	55,55,55,55	0
54	MG	DA	3045	1/1	0.92	0.59	8.48	39,39,39,39	0
54	MG	BA	3150	1/1	0.73	0.61	8.47	97,97,97,97	0
54	MG	DA	2932	1/1	0.98	0.24	8.35	24,24,24,24	0
54	MG	BA	3022	1/1	0.76	0.38	8.25	83,83,83,83	0
54	MG	DA	3059	1/1	0.98	0.34	8.21	39,39,39,39	0
54	MG	DA	3173	1/1	0.87	0.29	7.83	73,73,73,73	0
54	MG	DA	3106	1/1	0.90	0.26	7.78	69,69,69,69	0
54	MG	DA	3158	1/1	0.88	0.40	7.74	42,42,42,42	0
54	MG	BA	3068	1/1	0.90	0.25	7.69	55,55,55,55	0
54	MG	BA	3114	1/1	0.67	0.48	7.67	69,69,69,69	0
54	MG	CA	1891	1/1	0.71	0.43	7.65	113,113,113,113	0
54	MG	DA	3267	1/1	0.69	0.39	7.63	43,43,43,43	0
54	MG	DA	3414	1/1	0.94	0.32	7.60	55,55,55,55	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
54	MG	DA	2997	1/1	0.95	0.29	7.54	46,46,46,46	0
54	MG	DA	3074	1/1	0.95	0.32	7.51	43,43,43,43	0
54	MG	DA	2973	1/1	0.99	0.30	7.42	50,50,50,50	0
54	MG	BA	3064	1/1	0.89	0.37	7.36	61,61,61,61	0
54	MG	AA	1828	1/1	0.94	0.38	7.31	82,82,82,82	0
54	MG	DA	3258	1/1	0.96	0.28	7.21	67,67,67,67	0
54	MG	DA	3607	1/1	0.93	0.41	7.18	79,79,79,79	0
54	MG	BA	3004	1/1	0.95	0.28	7.14	49,49,49,49	0
54	MG	BA	3051	1/1	0.97	0.49	7.13	62,62,62,62	0
54	MG	CA	1737	1/1	0.86	0.47	7.12	70,70,70,70	0
54	MG	BA	2923	1/1	0.94	0.49	7.08	13,13,13,13	0
54	MG	DA	3064	1/1	0.96	0.24	6.97	127,127,127,127	0
54	MG	DA	3622	1/1	0.94	0.35	6.84	45,45,45,45	0
54	MG	BA	3187	1/1	0.62	0.75	6.81	73,73,73,73	0
54	MG	DA	3357	1/1	0.98	0.35	6.78	169,169,169,169	0
54	MG	DA	3311	1/1	0.70	0.55	6.70	95,95,95,95	0
54	MG	BA	2945	1/1	0.95	0.29	6.62	56,56,56,56	0
54	MG	BA	2912	1/1	0.89	0.50	6.57	19,19,19,19	0
54	MG	BA	3164	1/1	0.92	0.34	6.47	75,75,75,75	0
54	MG	BA	3442	1/1	0.96	0.51	6.38	26,26,26,26	0
54	MG	AA	1778	1/1	0.85	0.62	6.38	113,113,113,113	0
54	MG	DA	3840	1/1	0.88	0.38	6.37	81,81,81,81	0
54	MG	DA	3180	1/1	0.80	0.30	6.34	78,78,78,78	0
54	MG	BA	3243	1/1	0.80	0.51	6.29	85,85,85,85	0
54	MG	DA	2934	1/1	0.97	0.40	6.27	13,13,13,13	0
54	MG	CA	1733	1/1	0.94	0.56	6.16	78,78,78,78	0
54	MG	BA	2909	1/1	0.98	0.48	6.15	16,16,16,16	0
54	MG	AA	1632	1/1	0.89	0.34	6.08	65,65,65,65	0
54	MG	DA	3360	1/1	0.82	0.33	6.04	107,107,107,107	0
54	MG	BX	102	1/1	0.88	0.35	5.97	103,103,103,103	0
54	MG	BA	3443	1/1	0.98	0.34	5.97	36,36,36,36	0
54	MG	DA	3152	1/1	0.90	0.38	5.96	43,43,43,43	0
54	MG	BA	2919	1/1	0.98	0.32	5.86	14,14,14,14	0
54	MG	AA	1729	1/1	0.72	0.47	5.80	107,107,107,107	0
54	MG	BA	2955	1/1	0.97	0.32	5.80	63,63,63,63	0
54	MG	BA	2943	1/1	0.97	0.17	5.78	34,34,34,34	0
54	MG	BA	3434	1/1	0.97	0.27	5.75	19,19,19,19	0
54	MG	CA	1881	1/1	0.81	0.30	5.67	74,74,74,74	0
54	MG	DA	3056	1/1	0.97	0.24	5.67	25,25,25,25	0
54	MG	BA	3438	1/1	0.97	0.57	5.67	29,29,29,29	0
54	MG	DA	2904	1/1	0.98	0.30	5.63	16,16,16,16	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
54	MG	DA	2926	1/1	0.97	0.39	5.60	25,25,25,25	0
54	MG	DA	3048	1/1	0.87	0.29	5.50	51,51,51,51	0
54	MG	BB	218	1/1	0.98	0.36	5.47	79,79,79,79	0
54	MG	BA	3126	1/1	0.86	0.40	5.45	60,60,60,60	0
54	MG	AA	1665	1/1	0.87	0.32	5.45	70,70,70,70	0
54	MG	BA	3110	1/1	0.84	0.48	5.44	72,72,72,72	0
54	MG	AA	1625	1/1	0.94	0.22	5.38	53,53,53,53	0
54	MG	BA	3469	1/1	0.89	0.25	5.38	58,58,58,58	0
54	MG	DA	3005	1/1	0.98	0.36	5.35	41,41,41,41	0
54	MG	DD	301	1/1	0.91	0.50	5.31	39,39,39,39	0
54	MG	BA	3411	1/1	0.96	0.34	5.30	90,90,90,90	0
54	MG	DA	3121	1/1	0.95	0.35	5.23	38,38,38,38	0
54	MG	DA	2906	1/1	0.98	0.32	5.18	16,16,16,16	0
54	MG	DA	3343	1/1	0.78	0.30	5.16	97,97,97,97	0
54	MG	DA	3529	1/1	0.85	0.23	5.13	61,61,61,61	0
54	MG	BA	3673	1/1	0.80	0.28	5.08	92,92,92,92	0
54	MG	DA	3320	1/1	0.86	0.24	5.05	80,80,80,80	0
54	MG	DB	204	1/1	0.89	0.28	5.04	56,56,56,56	0
54	MG	BA	3169	1/1	0.84	0.26	4.99	43,43,43,43	0
54	MG	DA	3672	1/1	0.93	0.17	4.98	80,80,80,80	0
54	MG	BA	2914	1/1	0.96	0.45	4.93	24,24,24,24	0
54	MG	DA	2938	1/1	0.98	0.34	4.93	7,7,7,7	0
54	MG	DB	216	1/1	0.69	0.38	4.92	101,101,101,101	0
54	MG	BA	3575	1/1	0.95	0.38	4.90	59,59,59,59	0
54	MG	CA	1788	1/1	0.91	0.38	4.88	95,95,95,95	0
54	MG	BA	3038	1/1	0.74	0.28	4.81	63,63,63,63	0
54	MG	CA	1651	1/1	0.87	0.40	4.79	111,111,111,111	0
54	MG	DA	3014	1/1	0.91	0.25	4.74	52,52,52,52	0
54	MG	AA	1641	1/1	0.74	0.35	4.68	66,66,66,66	0
54	MG	BA	3327	1/1	0.85	0.79	4.65	86,86,86,86	0
54	MG	BE	302	1/1	0.93	0.44	4.60	93,93,93,93	0
54	MG	BA	3634	1/1	0.49	0.17	4.50	98,98,98,98	0
54	MG	DA	3197	1/1	0.92	0.17	4.47	46,46,46,46	0
54	MG	AA	1720	1/1	0.93	0.83	4.40	101,101,101,101	0
54	MG	DA	2980	1/1	0.98	0.37	4.32	12,12,12,12	0
54	MG	DA	3126	1/1	0.96	0.32	4.31	64,64,64,64	0
54	MG	DA	3851	1/1	0.94	0.27	4.23	74,74,74,74	0
54	MG	DA	3590	1/1	0.90	0.45	4.21	47,47,47,47	0
54	MG	AA	1609	1/1	0.98	0.33	4.18	26,26,26,26	0
54	MG	CA	1862	1/1	0.92	0.43	4.14	96,96,96,96	0
54	MG	AA	1635	1/1	0.99	0.27	4.07	37,37,37,37	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
54	MG	BA	3178	1/1	0.93	0.27	4.05	74,74,74,74	0
54	MG	DA	3638	1/1	0.95	0.31	4.04	22,22,22,22	0
54	MG	DA	3415	1/1	0.90	0.48	4.04	66,66,66,66	0
54	MG	BA	2990	1/1	0.94	0.43	4.03	35,35,35,35	0
54	MG	DA	3506	1/1	0.87	0.40	3.90	101,101,101,101	0
54	MG	BA	3289	1/1	0.96	0.39	3.90	61,61,61,61	0
54	MG	DA	3310	1/1	0.85	0.34	3.74	54,54,54,54	0
54	MG	DA	3108	1/1	0.88	0.26	3.74	54,54,54,54	0
54	MG	BA	3471	1/1	0.81	0.27	3.71	70,70,70,70	0
54	MG	BA	3098	1/1	0.95	0.27	3.69	90,90,90,90	0
54	MG	DA	3353	1/1	0.89	0.25	3.63	105,105,105,105	0
54	MG	AT	202	1/1	0.91	0.39	3.59	80,80,80,80	0
54	MG	BA	2991	1/1	0.97	0.22	3.41	42,42,42,42	0
54	MG	DA	3153	1/1	0.97	0.35	3.40	33,33,33,33	0
54	MG	BC	303	1/1	0.83	0.42	3.36	95,95,95,95	0
54	MG	DA	3455	1/1	0.94	0.90	3.34	97,97,97,97	0
54	MG	BA	3019	1/1	0.74	0.23	3.29	67,67,67,67	0
54	MG	DA	3052	1/1	0.97	0.27	3.11	50,50,50,50	0
54	MG	DA	3329	1/1	0.86	0.17	3.08	96,96,96,96	0
54	MG	BA	3238	1/1	0.92	0.44	3.06	49,49,49,49	0
54	MG	AA	1603	1/1	0.95	0.50	3.05	37,37,37,37	0
54	MG	BA	3047	1/1	0.93	0.23	3.05	72,72,72,72	0
54	MG	CA	1915	1/1	0.97	0.41	3.03	58,58,58,58	0
54	MG	BA	2975	1/1	0.98	0.26	3.00	58,58,58,58	0
54	MG	DA	2927	1/1	0.98	0.25	2.97	18,18,18,18	0
54	MG	AA	1870	1/1	0.93	0.25	2.91	98,98,98,98	0
54	MG	DA	3082	1/1	0.98	0.41	2.89	68,68,68,68	0
54	MG	DA	3736	1/1	0.95	0.23	2.87	134,134,134,134	0
54	MG	AD	303	1/1	0.94	0.47	2.87	68,68,68,68	0
54	MG	DA	3771	1/1	0.80	0.35	2.86	58,58,58,58	0
54	MG	DJ	205	1/1	0.88	0.50	2.76	91,91,91,91	0
54	MG	DW	101	1/1	0.85	0.33	2.75	56,56,56,56	0
54	MG	DA	3599	1/1	0.91	0.27	2.71	58,58,58,58	0
54	MG	BA	2981	1/1	0.96	0.26	2.70	56,56,56,56	0
54	MG	DA	3172	1/1	0.98	0.24	2.70	88,88,88,88	0
54	MG	DA	3268	1/1	0.92	0.25	2.67	51,51,51,51	0
54	MG	DA	3658	1/1	0.96	0.34	2.55	23,23,23,23	0
54	MG	DA	3289	1/1	0.91	0.26	2.55	44,44,44,44	0
54	MG	DA	3185	1/1	0.94	0.26	2.54	54,54,54,54	0
54	MG	BA	3334	1/1	0.96	0.24	2.52	60,60,60,60	0
54	MG	DA	2975	1/1	0.92	0.26	2.51	50,50,50,50	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
54	MG	CA	1760	1/1	0.81	0.27	2.48	66,66,66,66	0
54	MG	BA	2931	1/1	0.96	0.39	2.45	49,49,49,49	0
54	MG	DA	3098	1/1	0.96	0.24	2.44	46,46,46,46	0
54	MG	DA	3412	1/1	0.95	0.29	2.41	86,86,86,86	0
54	MG	AA	1630	1/1	0.81	0.27	2.39	81,81,81,81	0
54	MG	DX	101	1/1	0.85	0.37	2.38	103,103,103,103	0
54	MG	AA	1615	1/1	0.97	0.35	2.37	67,67,67,67	0
54	MG	BM	203	1/1	0.92	0.47	2.35	90,90,90,90	0
54	MG	BA	3013	1/1	0.96	0.40	2.29	21,21,21,21	0
54	MG	CA	1659	1/1	0.90	0.36	2.29	100,100,100,100	0
54	MG	CA	1688	1/1	0.54	0.26	2.24	80,80,80,80	0
54	MG	BA	3459	1/1	0.92	0.34	2.24	30,30,30,30	0
54	MG	DA	3606	1/1	0.89	0.26	2.22	68,68,68,68	0
54	MG	AA	1610	1/1	0.91	0.36	2.19	29,29,29,29	0
54	MG	DA	3605	1/1	0.83	0.27	2.17	103,103,103,103	0
54	MG	DA	3431	1/1	0.94	0.24	2.16	94,94,94,94	0
54	MG	BA	3301	1/1	0.78	0.38	2.15	58,58,58,58	0
54	MG	BA	3467	1/1	0.96	0.38	2.10	30,30,30,30	0
54	MG	DA	3028	1/1	0.95	0.35	2.07	58,58,58,58	0
54	MG	DA	2943	1/1	0.98	0.28	2.07	5,5,5,5	0
54	MG	DA	3010	1/1	0.90	0.23	2.06	72,72,72,72	0
54	MG	CA	1943	1/1	0.81	0.39	2.04	87,87,87,87	0
54	MG	DA	3015	1/1	0.94	0.26	2.03	33,33,33,33	0
54	MG	DA	3589	1/1	0.97	0.26	2.00	51,51,51,51	0
54	MG	BA	3356	1/1	0.96	0.28	1.99	33,33,33,33	0
54	MG	DA	3142	1/1	0.93	0.27	1.99	61,61,61,61	0
54	MG	DA	3750	1/1	0.91	0.45	1.99	87,87,87,87	0
54	MG	DA	3040	1/1	0.78	0.34	1.98	76,76,76,76	0
54	MG	BA	3252	1/1	0.85	0.28	1.96	60,60,60,60	0
54	MG	AA	1719	1/1	0.94	0.24	1.96	66,66,66,66	0
54	MG	BA	3175	1/1	0.80	0.43	1.94	71,71,71,71	0
54	MG	BA	2960	1/1	0.95	0.54	1.90	39,39,39,39	0
54	MG	CA	1675	1/1	0.93	0.24	1.90	74,74,74,74	0
54	MG	BA	2917	1/1	0.99	0.26	1.89	6,6,6,6	0
54	MG	AA	1601	1/1	0.96	0.24	1.89	35,35,35,35	0
54	MG	BB	220	1/1	0.83	0.33	1.84	75,75,75,75	0
54	MG	BA	3237	1/1	0.81	0.30	1.81	64,64,64,64	0
54	MG	BA	3030	1/1	0.89	0.18	1.79	61,61,61,61	0
54	MG	BA	3515	1/1	0.82	0.38	1.77	54,54,54,54	0
54	MG	DA	3361	1/1	0.95	0.12	1.74	69,69,69,69	0
54	MG	DA	3410	1/1	0.81	0.32	1.72	48,48,48,48	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
54	MG	CN	102	1/1	0.91	1.05	1.70	134,134,134,134	0
54	MG	DA	3250	1/1	0.65	0.27	1.67	70,70,70,70	0
54	MG	BA	3256	1/1	0.75	0.17	1.66	105,105,105,105	0
54	MG	DS	201	1/1	0.84	0.37	1.57	98,98,98,98	0
54	MG	DA	3554	1/1	0.84	0.30	1.50	136,136,136,136	0
54	MG	DA	3146	1/1	0.96	0.22	1.47	31,31,31,31	0
54	MG	CA	1944	1/1	0.88	0.32	1.43	66,66,66,66	0
54	MG	DA	3186	1/1	0.89	0.41	1.40	82,82,82,82	0
54	MG	BM	202	1/1	0.95	0.30	1.39	85,85,85,85	0
54	MG	DA	3514	1/1	0.89	0.23	1.34	68,68,68,68	0
54	MG	CA	1921	1/1	0.87	0.21	1.25	97,97,97,97	0
54	MG	DA	3307	1/1	0.93	0.21	1.21	84,84,84,84	0
54	MG	DA	3176	1/1	0.94	0.31	1.20	50,50,50,50	0
54	MG	DA	3283	1/1	0.88	0.27	1.19	71,71,71,71	0
54	MG	AA	1614	1/1	0.81	0.20	1.09	68,68,68,68	0
54	MG	CA	1781	1/1	0.93	0.22	1.08	76,76,76,76	0
54	MG	BA	3430	1/1	0.97	0.27	1.07	6,6,6,6	0
54	MG	CA	1826	1/1	0.86	0.31	1.02	97,97,97,97	0
54	MG	BA	2948	1/1	0.90	0.27	0.99	62,62,62,62	0
54	MG	BA	3677	1/1	0.87	0.26	0.98	72,72,72,72	0
54	MG	DA	3625	1/1	0.94	0.32	0.97	84,84,84,84	0
54	MG	BA	3574	1/1	0.95	0.18	0.89	62,62,62,62	0
54	MG	DA	3632	1/1	0.73	0.29	0.88	111,111,111,111	0
54	MG	DA	3044	1/1	0.90	0.19	0.83	45,45,45,45	0
54	MG	DA	3099	1/1	0.92	0.18	0.81	49,49,49,49	0
54	MG	AA	1784	1/1	0.72	0.42	0.80	89,89,89,89	0
54	MG	CA	1762	1/1	0.96	0.21	0.76	114,114,114,114	0
54	MG	CA	1607	1/1	0.93	0.18	0.75	62,62,62,62	0
54	MG	DA	2990	1/1	0.95	0.21	0.74	20,20,20,20	0
54	MG	DA	3037	1/1	0.93	0.18	0.74	44,44,44,44	0
54	MG	DA	3484	1/1	0.96	0.25	0.72	21,21,21,21	0
54	MG	B5	101	1/1	0.91	0.47	0.68	78,78,78,78	0
54	MG	DA	3125	1/1	0.96	0.20	0.66	36,36,36,36	0
54	MG	BA	3107	1/1	0.99	0.31	0.66	25,25,25,25	0
54	MG	DB	215	1/1	0.86	0.28	0.60	103,103,103,103	0
54	MG	CM	201	1/1	0.92	0.89	0.59	85,85,85,85	0
54	MG	DD	305	1/1	0.76	0.41	0.57	52,52,52,52	0
54	MG	DA	3641	1/1	0.95	0.25	0.55	13,13,13,13	0
54	MG	AE	202	1/1	0.85	0.21	0.50	173,173,173,173	0
54	MG	DJ	204	1/1	0.93	0.50	0.48	32,32,32,32	0
54	MG	CA	1702	1/1	0.96	0.24	0.47	114,114,114,114	0
54	MG	BA	2933	1/1	0.95	0.24	0.45	36,36,36,36	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
54	MG	BA	3307	1/1	0.94	0.19	0.43	63,63,63,63	0
54	MG	D4	102	1/1	0.88	0.23	0.42	65,65,65,65	0
54	MG	DA	3104	1/1	0.89	0.24	0.41	49,49,49,49	0
54	MG	DA	3680	1/1	0.96	0.20	0.41	41,41,41,41	0
54	MG	DA	3050	1/1	0.98	0.21	0.35	89,89,89,89	0
54	MG	CA	1950	1/1	0.77	0.36	0.35	107,107,107,107	0
54	MG	BA	2942	1/1	0.88	0.21	0.35	19,19,19,19	0
54	MG	CA	1690	1/1	0.86	0.16	0.32	65,65,65,65	0
54	MG	BA	3403	1/1	0.86	0.26	0.31	106,106,106,106	0
54	MG	DA	2972	1/1	0.97	0.19	0.29	37,37,37,37	0
54	MG	DA	3137	1/1	0.94	0.29	0.28	29,29,29,29	0
54	MG	DA	3613	1/1	0.55	0.32	0.27	121,121,121,121	0
54	MG	BA	3547	1/1	0.96	0.20	0.26	87,87,87,87	0
54	MG	BJ	202	1/1	0.94	0.31	0.24	69,69,69,69	0
54	MG	CD	304	1/1	0.89	0.50	0.22	108,108,108,108	0
54	MG	AA	1649	1/1	0.77	0.16	0.21	74,74,74,74	0
54	MG	CA	1657	1/1	0.79	0.17	0.20	60,60,60,60	0
54	MG	DA	2999	1/1	0.93	0.15	0.19	69,69,69,69	0
54	MG	DA	2995	1/1	0.95	0.21	0.17	26,26,26,26	0
54	MG	DA	2984	1/1	0.97	0.22	0.17	22,22,22,22	0
54	MG	BA	2988	1/1	0.97	0.21	0.15	40,40,40,40	0
54	MG	BA	3546	1/1	0.96	0.20	0.15	67,67,67,67	0
54	MG	BA	3370	1/1	0.79	0.27	0.13	76,76,76,76	0
54	MG	BA	3615	1/1	0.93	0.18	0.11	93,93,93,93	0
54	MG	DA	3634	1/1	0.84	1.08	0.10	106,106,106,106	0
54	MG	BA	2936	1/1	0.95	0.18	0.06	32,32,32,32	0
54	MG	DB	212	1/1	0.92	0.24	0.05	85,85,85,85	0
54	MG	DA	3001	1/1	0.89	0.18	0.03	25,25,25,25	0
54	MG	BA	3562	1/1	0.91	0.20	-0.01	40,40,40,40	0
54	MG	AA	1841	1/1	0.95	0.17	-0.01	83,83,83,83	0
54	MG	BA	2973	1/1	0.82	0.15	-0.02	60,60,60,60	0
54	MG	DA	3326	1/1	0.94	0.32	-0.04	81,81,81,81	0
54	MG	CA	1764	1/1	0.75	0.15	-0.16	65,65,65,65	0
54	MG	BA	2989	1/1	0.94	0.20	-0.16	50,50,50,50	0
55	ZN	AD	301	1/1	0.98	0.29	-0.21	51,51,51,51	0
54	MG	BW	101	1/1	0.91	0.20	-0.22	57,57,57,57	0
54	MG	AA	1658	1/1	0.81	0.18	-0.23	84,84,84,84	0
54	MG	BA	3522	1/1	0.89	0.27	-0.23	54,54,54,54	0
54	MG	AA	1835	1/1	0.94	0.23	-0.26	158,158,158,158	0
54	MG	DQ	201	1/1	0.88	0.23	-0.27	57,57,57,57	0
54	MG	CA	1811	1/1	-0.03	0.28	-0.28	181,181,181,181	0
54	MG	BA	3131	1/1	0.92	0.15	-0.28	47,47,47,47	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
54	MG	DL	202	1/1	0.64	0.23	-0.32	146,146,146,146	0
54	MG	CA	1620	1/1	0.94	0.20	-0.33	51,51,51,51	0
54	MG	AA	1660	1/1	0.87	0.17	-0.37	122,122,122,122	0
54	MG	DA	3338	1/1	0.98	0.14	-0.38	123,123,123,123	0
54	MG	BA	3482	1/1	0.95	0.21	-0.39	40,40,40,40	0
54	MG	AA	1624	1/1	0.96	0.18	-0.40	58,58,58,58	0
54	MG	BX	101	1/1	0.91	0.21	-0.41	71,71,71,71	0
54	MG	DK	201	1/1	0.86	0.16	-0.48	87,87,87,87	0
54	MG	BA	3056	1/1	0.94	0.19	-0.49	31,31,31,31	0
54	MG	CA	1854	1/1	0.87	0.14	-0.50	111,111,111,111	0
54	MG	BA	3099	1/1	0.93	0.11	-0.52	56,56,56,56	0
54	MG	CK	201	1/1	0.78	0.17	-0.60	136,136,136,136	0
54	MG	AB	301	1/1	0.94	0.30	-0.61	92,92,92,92	0
54	MG	DA	2978	1/1	0.94	0.24	-0.62	23,23,23,23	0
54	MG	BA	3125	1/1	0.91	0.15	-0.63	78,78,78,78	0
54	MG	AM	201	1/1	0.76	0.23	-0.69	126,126,126,126	0
54	MG	BA	3535	1/1	0.95	0.21	-0.70	95,95,95,95	0
54	MG	DA	3281	1/1	0.96	0.16	-0.71	62,62,62,62	0
54	MG	CA	1710	1/1	0.79	0.16	-0.72	169,169,169,169	0
54	MG	AB	302	1/1	0.96	0.24	-0.74	57,57,57,57	0
54	MG	CA	1689	1/1	0.91	0.14	-0.74	130,130,130,130	0
54	MG	DA	3016	1/1	0.98	0.18	-0.76	47,47,47,47	0
54	MG	BA	3104	1/1	0.93	0.15	-0.77	49,49,49,49	0
54	MG	DI	101	1/1	0.79	0.29	-0.79	104,104,104,104	0
54	MG	DP	201	1/1	0.92	0.26	-0.79	72,72,72,72	0
55	ZN	CN	101	1/1	0.83	0.20	-0.82	244,244,244,244	0
54	MG	DA	3023	1/1	0.98	0.15	-0.83	23,23,23,23	0
54	MG	DA	3214	1/1	0.99	0.17	-0.85	34,34,34,34	0
54	MG	CA	1703	1/1	0.81	0.21	-0.95	113,113,113,113	0
54	MG	BA	3241	1/1	0.95	0.13	-0.99	69,69,69,69	0
54	MG	BA	3581	1/1	0.99	0.14	-1.01	127,127,127,127	0
54	MG	DA	3796	1/1	0.95	0.15	-1.03	67,67,67,67	0
54	MG	DA	3000	1/1	0.92	0.20	-1.03	23,23,23,23	0
54	MG	DA	2941	1/1	0.90	0.18	-1.05	21,21,21,21	0
54	MG	AA	1701	1/1	0.82	0.19	-1.13	84,84,84,84	0
54	MG	BA	2980	1/1	0.96	0.18	-1.15	51,51,51,51	0
54	MG	DA	3306	1/1	0.83	0.15	-1.19	152,152,152,152	0
54	MG	DA	3119	1/1	0.96	0.15	-1.20	74,74,74,74	0
54	MG	BA	3123	1/1	0.90	0.14	-1.21	40,40,40,40	0
55	ZN	AN	101	1/1	0.95	0.13	-1.21	144,144,144,144	0
54	MG	B2	101	1/1	0.93	0.11	-1.22	27,27,27,27	0
54	MG	CA	1747	1/1	0.90	0.17	-1.26	69,69,69,69	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
54	MG	AE	201	1/1	0.86	0.11	-1.27	123,123,123,123	0
54	MG	AA	1802	1/1	0.91	0.09	-1.33	62,62,62,62	0
54	MG	CA	1860	1/1	0.94	0.21	-1.33	68,68,68,68	0
54	MG	AA	1644	1/1	0.92	0.14	-1.33	62,62,62,62	0
54	MG	BA	3014	1/1	0.92	0.08	-1.36	45,45,45,45	0
54	MG	BA	2982	1/1	0.94	0.12	-1.37	38,38,38,38	0
54	MG	BA	2950	1/1	0.98	0.16	-1.37	68,68,68,68	0
54	MG	CA	1616	1/1	0.96	0.14	-1.48	94,94,94,94	0
54	MG	DE	301	1/1	0.86	0.17	-1.50	32,32,32,32	0
54	MG	DA	2963	1/1	0.87	0.10	-1.61	32,32,32,32	0
54	MG	DB	217	1/1	0.95	0.09	-1.62	93,93,93,93	0
54	MG	CA	1720	1/1	0.90	0.22	-1.65	69,69,69,69	0
54	MG	CA	1937	1/1	0.96	0.27	-1.69	100,100,100,100	0
54	MG	CA	1749	1/1	0.98	0.11	-1.70	77,77,77,77	0
54	MG	BE	301	1/1	0.77	0.13	-1.82	75,75,75,75	0
54	MG	AS	101	1/1	0.40	0.12	-1.84	106,106,106,106	0
54	MG	BA	3477	1/1	0.96	0.13	-1.88	50,50,50,50	0
54	MG	DA	3117	1/1	0.96	0.12	-1.93	31,31,31,31	0
54	MG	DX	102	1/1	0.92	0.12	-1.96	91,91,91,91	0
54	MG	BA	3412	1/1	0.98	0.13	-1.96	49,49,49,49	0
54	MG	AD	302	1/1	0.85	0.18	-1.97	76,76,76,76	0
54	MG	DA	3139	1/1	0.97	0.15	-2.05	40,40,40,40	0
54	MG	CC	301	1/1	0.78	0.09	-2.05	93,93,93,93	0
54	MG	DB	219	1/1	0.89	0.17	-2.16	73,73,73,73	0
54	MG	DA	3591	1/1	0.95	0.10	-2.21	90,90,90,90	0
54	MG	AA	1684	1/1	0.91	0.11	-2.26	71,71,71,71	0
54	MG	DA	2948	1/1	0.99	0.13	-2.50	21,21,21,21	0
55	ZN	CD	301	1/1	0.94	0.28	-2.51	116,116,116,116	0
54	MG	DA	3120	1/1	0.87	0.10	-2.62	61,61,61,61	0
54	MG	DA	3242	1/1	0.91	0.11	-2.65	59,59,59,59	0
54	MG	BA	3080	1/1	0.94	0.11	-2.69	43,43,43,43	0
54	MG	BA	3387	1/1	0.96	0.13	-2.77	84,84,84,84	0
54	MG	BA	2966	1/1	0.98	0.14	-2.81	38,38,38,38	0
54	MG	BA	3154	1/1	0.94	0.08	-3.45	139,139,139,139	0
54	MG	AA	1634	1/1	0.97	0.10	-3.47	48,48,48,48	0
54	MG	DA	3130	1/1	0.96	0.08	-3.51	68,68,68,68	0
54	MG	DA	3184	1/1	0.93	0.12	-3.67	38,38,38,38	0
54	MG	CA	1770	1/1	0.80	0.09	-4.19	125,125,125,125	0
54	MG	D2	102	1/1	0.90	0.08	-4.21	42,42,42,42	0
54	MG	BA	3015	1/1	0.97	0.10	-7.40	55,55,55,55	0
54	MG	DA	3301	1/1	0.92	0.68	-	87,87,87,87	0
54	MG	DA	3279	1/1	0.89	0.37	-	38,38,38,38	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
54	MG	DA	3422	1/1	0.72	0.62	-	62,62,62,62	0
54	MG	DA	3195	1/1	0.88	0.25	-	83,83,83,83	0
54	MG	DB	227	1/1	0.91	0.16	-	104,104,104,104	0
54	MG	BA	3505	1/1	0.80	0.29	-	81,81,81,81	0
54	MG	DA	3224	1/1	0.81	0.40	-	55,55,55,55	0
54	MG	BA	3589	1/1	0.34	0.19	-	105,105,105,105	0
54	MG	DA	3461	1/1	0.82	0.38	-	88,88,88,88	0
54	MG	BA	3408	1/1	0.93	0.51	-	86,86,86,86	0
54	MG	DV	203	1/1	0.92	0.15	-	128,128,128,128	0
54	MG	CA	1755	1/1	0.95	0.45	-	69,69,69,69	0
54	MG	DA	3649	1/1	0.92	1.06	-	50,50,50,50	0
54	MG	BA	3151	1/1	0.88	0.45	-	70,70,70,70	0
54	MG	BA	3517	1/1	0.94	0.42	-	87,87,87,87	0
54	MG	CA	1894	1/1	0.88	0.47	-	97,97,97,97	0
54	MG	AK	202	1/1	0.77	0.10	-	101,101,101,101	0
54	MG	BA	2956	1/1	0.85	0.63	-	49,49,49,49	0
54	MG	DA	3748	1/1	0.88	0.24	-	61,61,61,61	0
54	MG	DA	3707	1/1	0.85	0.20	-	54,54,54,54	0
54	MG	CA	1784	1/1	0.90	0.24	-	94,94,94,94	0
54	MG	BA	3250	1/1	0.86	0.13	-	67,67,67,67	0
54	MG	CA	1857	1/1	0.95	0.33	-	76,76,76,76	0
54	MG	DA	3490	1/1	0.94	0.21	-	70,70,70,70	0
54	MG	DA	3252	1/1	0.92	0.13	-	52,52,52,52	0
54	MG	CA	1869	1/1	0.93	0.08	-	96,96,96,96	0
54	MG	BA	2977	1/1	0.97	0.40	-	69,69,69,69	0
54	MG	DA	3546	1/1	0.91	0.20	-	56,56,56,56	0
54	MG	BA	3468	1/1	0.79	0.55	-	57,57,57,57	0
54	MG	DA	3492	1/1	0.92	0.18	-	102,102,102,102	0
54	MG	D5	103	1/1	0.90	0.38	-	71,71,71,71	0
54	MG	BA	2949	1/1	0.85	0.21	-	65,65,65,65	0
54	MG	BA	3364	1/1	0.90	0.36	-	107,107,107,107	0
54	MG	DA	3619	1/1	0.85	0.95	-	83,83,83,83	0
54	MG	DA	3566	1/1	0.82	0.29	-	91,91,91,91	0
54	MG	CA	1630	1/1	0.97	0.25	-	49,49,49,49	0
54	MG	DA	3442	1/1	0.89	0.63	-	71,71,71,71	0
54	MG	DA	3300	1/1	0.95	0.23	-	46,46,46,46	0
54	MG	BA	3044	1/1	0.87	0.26	-	39,39,39,39	0
54	MG	CA	1906	1/1	0.82	0.32	-	80,80,80,80	0
54	MG	AA	1821	1/1	0.87	0.51	-	91,91,91,91	0
54	MG	DA	3820	1/1	0.90	0.94	-	59,59,59,59	0
54	MG	BA	3319	1/1	0.95	0.16	-	79,79,79,79	0
54	MG	DA	3365	1/1	0.85	0.34	-	83,83,83,83	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
54	MG	AA	1852	1/1	0.92	0.78	-	84,84,84,84	0
54	MG	CA	1832	1/1	0.95	0.11	-	92,92,92,92	0
54	MG	CA	1821	1/1	0.93	0.50	-	96,96,96,96	0
54	MG	BA	2915	1/1	0.92	0.34	-	20,20,20,20	0
54	MG	CA	1748	1/1	0.88	0.21	-	115,115,115,115	0
54	MG	CA	1609	1/1	0.96	0.25	-	66,66,66,66	0
54	MG	DA	3702	1/1	0.72	0.60	-	74,74,74,74	0
54	MG	BA	3413	1/1	0.76	0.41	-	65,65,65,65	0
54	MG	AA	1775	1/1	0.96	0.11	-	105,105,105,105	0
54	MG	BA	3053	1/1	0.90	0.45	-	69,69,69,69	0
54	MG	BA	3065	1/1	0.95	0.22	-	59,59,59,59	0
54	MG	DA	3149	1/1	0.94	0.25	-	60,60,60,60	0
54	MG	DA	2924	1/1	0.80	0.51	-	68,68,68,68	0
54	MG	B2	103	1/1	0.84	0.09	-	75,75,75,75	0
54	MG	BA	3541	1/1	0.81	0.69	-	89,89,89,89	0
54	MG	BA	3519	1/1	0.91	0.15	-	50,50,50,50	0
54	MG	DA	3731	1/1	0.95	0.14	-	60,60,60,60	0
54	MG	DA	3550	1/1	0.91	0.14	-	65,65,65,65	0
54	MG	BA	3127	1/1	0.88	0.77	-	61,61,61,61	0
54	MG	DA	3577	1/1	0.92	0.40	-	73,73,73,73	0
54	MG	DA	3089	1/1	0.96	0.40	-	71,71,71,71	0
54	MG	AV	6207	1/1	0.91	0.11	-	78,78,78,78	0
54	MG	BA	3623	1/1	0.67	0.65	-	64,64,64,64	0
54	MG	CA	1729	1/1	0.91	0.16	-	58,58,58,58	0
54	MG	CA	1899	1/1	0.72	0.10	-	98,98,98,98	0
54	MG	BA	2928	1/1	0.95	0.51	-	28,28,28,28	0
54	MG	CQ	202	1/1	0.90	0.20	-	79,79,79,79	0
54	MG	DA	3664	1/1	0.56	1.62	-	109,109,109,109	0
54	MG	DA	3332	1/1	0.66	0.79	-	88,88,88,88	0
54	MG	CA	1647	1/1	0.82	0.27	-	94,94,94,94	0
54	MG	DA	2921	1/1	0.95	0.13	-	32,32,32,32	0
54	MG	CA	1797	1/1	0.70	0.27	-	133,133,133,133	0
54	MG	DA	3653	1/1	0.92	0.52	-	51,51,51,51	0
54	MG	BA	3233	1/1	0.98	0.61	-	97,97,97,97	0
54	MG	CA	1823	1/1	0.96	0.12	-	61,61,61,61	0
54	MG	CA	1671	1/1	0.98	0.26	-	51,51,51,51	0
54	MG	BA	3426	1/1	0.77	0.35	-	87,87,87,87	0
54	MG	BA	3369	1/1	0.95	0.30	-	108,108,108,108	0
54	MG	DA	2976	1/1	0.78	1.34	-	68,68,68,68	0
54	MG	CF	201	1/1	0.65	0.24	-	131,131,131,131	0
54	MG	BA	3166	1/1	0.75	0.90	-	86,86,86,86	0
54	MG	BA	3460	1/1	0.93	0.22	-	47,47,47,47	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
54	MG	CA	1642	1/1	0.95	0.18	-	79,79,79,79	0
54	MG	DA	3165	1/1	0.97	0.11	-	59,59,59,59	0
54	MG	DB	225	1/1	0.78	0.18	-	72,72,72,72	0
54	MG	AA	1748	1/1	0.86	0.21	-	56,56,56,56	0
54	MG	DA	3500	1/1	0.88	0.23	-	108,108,108,108	0
54	MG	AA	1779	1/1	0.90	0.71	-	108,108,108,108	0
54	MG	CA	1715	1/1	0.97	0.42	-	107,107,107,107	0
54	MG	DP	202	1/1	0.70	0.10	-	90,90,90,90	0
54	MG	BA	3205	1/1	0.83	0.48	-	70,70,70,70	0
54	MG	DA	3437	1/1	0.81	0.66	-	120,120,120,120	0
54	MG	DA	3374	1/1	0.99	0.14	-	80,80,80,80	0
54	MG	DA	3218	1/1	0.94	0.11	-	79,79,79,79	0
54	MG	CA	1684	1/1	0.93	0.49	-	63,63,63,63	0
54	MG	BA	3039	1/1	0.93	0.14	-	54,54,54,54	0
54	MG	BA	3636	1/1	0.93	0.77	-	76,76,76,76	0
54	MG	DA	3787	1/1	0.72	1.21	-	114,114,114,114	0
54	MG	BA	3320	1/1	0.93	0.41	-	71,71,71,71	0
54	MG	CA	1752	1/1	0.92	0.33	-	71,71,71,71	0
54	MG	DA	3156	1/1	0.90	0.21	-	52,52,52,52	0
54	MG	CA	1707	1/1	0.95	0.14	-	102,102,102,102	0
54	MG	BA	3341	1/1	0.12	0.66	-	125,125,125,125	0
54	MG	BA	3218	1/1	0.89	0.24	-	55,55,55,55	0
54	MG	DA	3439	1/1	0.94	0.27	-	52,52,52,52	0
54	MG	DA	3369	1/1	0.99	0.10	-	99,99,99,99	0
54	MG	AA	1810	1/1	0.96	0.24	-	86,86,86,86	0
54	MG	BA	3026	1/1	0.79	0.15	-	68,68,68,68	0
54	MG	BA	3216	1/1	0.92	0.18	-	87,87,87,87	0
54	MG	DA	3244	1/1	0.84	0.55	-	85,85,85,85	0
54	MG	DA	3836	1/1	0.20	0.63	-	135,135,135,135	0
54	MG	AA	1607	1/1	0.81	0.60	-	46,46,46,46	0
54	MG	CA	1683	1/1	0.84	0.34	-	72,72,72,72	0
54	MG	BA	3470	1/1	0.86	0.30	-	71,71,71,71	0
54	MG	DA	3384	1/1	0.83	0.30	-	84,84,84,84	0
54	MG	DA	3179	1/1	0.62	0.57	-	105,105,105,105	0
54	MG	DJ	206	1/1	0.88	0.14	-	135,135,135,135	0
54	MG	DA	3303	1/1	0.96	0.36	-	67,67,67,67	0
54	MG	DA	2945	1/1	0.94	0.58	-	29,29,29,29	0
54	MG	CA	1816	1/1	0.82	0.41	-	107,107,107,107	0
54	MG	AA	1811	1/1	0.93	0.19	-	61,61,61,61	0
54	MG	CA	1952	1/1	0.81	0.29	-	88,88,88,88	0
54	MG	DA	3521	1/1	0.93	0.24	-	101,101,101,101	0
54	MG	BA	3305	1/1	0.89	0.10	-	87,87,87,87	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
54	MG	DA	3849	1/1	0.88	0.17	-	89,89,89,89	0
54	MG	CA	1901	1/1	0.89	0.30	-	76,76,76,76	0
54	MG	DA	3069	1/1	0.79	0.46	-	61,61,61,61	0
54	MG	CA	1837	1/1	0.75	0.61	-	83,83,83,83	0
54	MG	DV	204	1/1	0.85	0.28	-	96,96,96,96	0
54	MG	DA	3071	1/1	0.99	0.16	-	49,49,49,49	0
54	MG	BA	3016	1/1	0.95	0.50	-	54,54,54,54	0
54	MG	BA	3595	1/1	0.83	0.34	-	108,108,108,108	0
54	MG	DA	3295	1/1	0.77	0.40	-	107,107,107,107	0
54	MG	BA	3423	1/1	0.95	0.39	-	72,72,72,72	0
54	MG	CA	1819	1/1	0.83	0.26	-	109,109,109,109	0
54	MG	DA	3525	1/1	0.95	0.75	-	73,73,73,73	0
54	MG	DD	303	1/1	0.72	0.17	-	70,70,70,70	0
54	MG	BA	3496	1/1	0.82	0.16	-	67,67,67,67	0
54	MG	DA	3123	1/1	0.93	0.42	-	59,59,59,59	0
54	MG	DA	3597	1/1	0.95	0.13	-	124,124,124,124	0
54	MG	BA	3361	1/1	0.83	0.27	-	154,154,154,154	0
54	MG	CA	1709	1/1	0.95	0.05	-	107,107,107,107	0
54	MG	DA	2959	1/1	0.92	0.30	-	42,42,42,42	0
54	MG	DA	3196	1/1	0.96	0.20	-	84,84,84,84	0
54	MG	BA	3100	1/1	0.94	0.31	-	79,79,79,79	0
54	MG	BA	3322	1/1	0.79	0.36	-	71,71,71,71	0
54	MG	DA	3533	1/1	0.88	0.25	-	60,60,60,60	0
54	MG	DA	3710	1/1	0.97	0.15	-	92,92,92,92	0
54	MG	DA	3386	1/1	0.93	0.13	-	76,76,76,76	0
54	MG	CA	1829	1/1	0.95	0.13	-	89,89,89,89	0
54	MG	DA	3094	1/1	0.89	0.41	-	67,67,67,67	0
54	MG	DA	3174	1/1	0.94	0.18	-	66,66,66,66	0
54	MG	AA	1674	1/1	0.82	0.95	-	113,113,113,113	0
54	MG	AG	202	1/1	0.82	0.31	-	129,129,129,129	0
54	MG	DA	3535	1/1	0.81	0.31	-	107,107,107,107	0
54	MG	AA	1663	1/1	0.90	0.22	-	54,54,54,54	0
54	MG	BA	3456	1/1	0.95	0.38	-	64,64,64,64	0
54	MG	BA	3464	1/1	0.70	0.73	-	67,67,67,67	0
54	MG	CA	1790	1/1	0.80	0.28	-	132,132,132,132	0
54	MG	DA	3743	1/1	0.71	0.27	-	105,105,105,105	0
54	MG	DA	3636	1/1	0.98	0.42	-	12,12,12,12	0
54	MG	BA	3475	1/1	0.77	1.13	-	67,67,67,67	0
54	MG	DA	3257	1/1	0.94	0.15	-	48,48,48,48	0
54	MG	DA	3447	1/1	0.92	0.25	-	73,73,73,73	0
54	MG	DB	205	1/1	0.94	0.32	-	78,78,78,78	0
54	MG	BA	3417	1/1	0.93	0.18	-	65,65,65,65	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
54	MG	BA	2946	1/1	0.97	0.21	-	37,37,37,37	0
54	MG	DA	3479	1/1	0.84	0.20	-	130,130,130,130	0
54	MG	CA	1807	1/1	0.73	0.37	-	125,125,125,125	0
54	MG	AV	6205	1/1	0.92	0.12	-	107,107,107,107	0
54	MG	DA	3436	1/1	0.72	1.02	-	101,101,101,101	0
54	MG	BA	3057	1/1	0.99	0.14	-	64,64,64,64	0
54	MG	BP	202	1/1	0.89	0.13	-	84,84,84,84	0
54	MG	AA	1670	1/1	0.57	0.14	-	93,93,93,93	0
54	MG	CA	1670	1/1	0.94	0.36	-	106,106,106,106	0
54	MG	CA	1672	1/1	0.95	0.14	-	137,137,137,137	0
54	MG	CA	1613	1/1	0.89	0.62	-	68,68,68,68	0
54	MG	AO	101	1/1	0.77	0.09	-	115,115,115,115	0
54	MG	CA	1835	1/1	0.93	0.23	-	76,76,76,76	0
54	MG	BA	3106	1/1	-0.09	0.22	-	142,142,142,142	0
54	MG	BA	2908	1/1	0.95	0.44	-	26,26,26,26	0
54	MG	DA	3434	1/1	0.91	0.27	-	89,89,89,89	0
54	MG	DA	3097	1/1	0.98	0.30	-	55,55,55,55	0
54	MG	CA	1961	1/1	0.98	0.12	-	63,63,63,63	0
54	MG	DV	202	1/1	0.72	0.34	-	115,115,115,115	0
54	MG	AA	1627	1/1	0.93	0.60	-	59,59,59,59	0
54	MG	BA	3630	1/1	0.90	0.37	-	151,151,151,151	0
54	MG	BA	3582	1/1	0.89	0.30	-	102,102,102,102	0
54	MG	CA	1624	1/1	0.95	0.43	-	94,94,94,94	0
54	MG	CA	1697	1/1	0.89	0.08	-	83,83,83,83	0
54	MG	AA	1680	1/1	0.57	0.42	-	140,140,140,140	0
54	MG	CA	1718	1/1	0.95	0.27	-	109,109,109,109	0
54	MG	AA	1650	1/1	0.99	0.22	-	92,92,92,92	0
54	MG	BA	3678	1/1	0.93	0.53	-	122,122,122,122	0
54	MG	DB	214	1/1	0.96	0.21	-	72,72,72,72	0
54	MG	AA	1739	1/1	0.91	0.27	-	81,81,81,81	0
54	MG	BB	216	1/1	0.54	0.69	-	154,154,154,154	0
54	MG	BA	3404	1/1	0.88	0.21	-	80,80,80,80	0
54	MG	BA	3362	1/1	0.76	0.93	-	130,130,130,130	0
54	MG	DA	3693	1/1	0.87	0.31	-	63,63,63,63	0
54	MG	DA	3383	1/1	0.77	0.34	-	112,112,112,112	0
54	MG	AA	1669	1/1	0.94	0.39	-	54,54,54,54	0
54	MG	AA	1611	1/1	0.88	0.45	-	55,55,55,55	0
54	MG	BA	3148	1/1	0.79	0.28	-	62,62,62,62	0
54	MG	BA	3349	1/1	0.68	0.33	-	95,95,95,95	0
54	MG	DA	3055	1/1	0.93	0.34	-	59,59,59,59	0
54	MG	BA	3399	1/1	0.84	0.38	-	104,104,104,104	0
54	MG	BA	3232	1/1	0.94	0.27	-	68,68,68,68	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
54	MG	CA	1628	1/1	0.89	0.63	-	84,84,84,84	0
54	MG	DA	3493	1/1	0.95	0.39	-	59,59,59,59	0
54	MG	DB	203	1/1	0.87	0.16	-	69,69,69,69	0
54	MG	BA	3001	1/1	0.85	0.37	-	65,65,65,65	0
54	MG	DA	3349	1/1	0.92	0.17	-	116,116,116,116	0
54	MG	BA	3197	1/1	0.74	0.18	-	84,84,84,84	0
54	MG	CA	1809	1/1	0.96	0.21	-	92,92,92,92	0
54	MG	BA	3018	1/1	0.62	0.84	-	88,88,88,88	0
54	MG	DA	3265	1/1	0.91	0.67	-	138,138,138,138	0
54	MG	DA	3340	1/1	0.71	0.33	-	59,59,59,59	0
54	MG	DA	3705	1/1	0.86	0.20	-	86,86,86,86	0
54	MG	DQ	202	1/1	0.82	0.20	-	97,97,97,97	0
54	MG	DA	3800	1/1	0.80	0.36	-	95,95,95,95	0
54	MG	DA	3854	1/1	0.87	0.10	-	124,124,124,124	0
54	MG	DA	3312	1/1	0.68	0.16	-	130,130,130,130	0
54	MG	DA	3448	1/1	0.96	0.17	-	99,99,99,99	0
54	MG	DB	224	1/1	0.97	0.26	-	95,95,95,95	0
54	MG	AA	1662	1/1	0.90	0.56	-	78,78,78,78	0
54	MG	CA	1822	1/1	0.90	0.38	-	93,93,93,93	0
54	MG	BA	3045	1/1	0.64	0.38	-	94,94,94,94	0
54	MG	BA	3651	1/1	0.84	0.49	-	77,77,77,77	0
54	MG	BA	2967	1/1	0.92	0.41	-	51,51,51,51	0
54	MG	DA	3560	1/1	0.85	0.44	-	83,83,83,83	0
54	MG	BA	3529	1/1	0.86	0.27	-	79,79,79,79	0
54	MG	AA	1721	1/1	0.95	0.32	-	56,56,56,56	0
54	MG	DA	3726	1/1	0.74	0.29	-	106,106,106,106	0
54	MG	DA	3392	1/1	0.90	0.50	-	76,76,76,76	0
54	MG	DA	3537	1/1	0.70	0.43	-	101,101,101,101	0
54	MG	DA	3665	1/1	0.47	0.87	-	137,137,137,137	0
54	MG	DA	3395	1/1	0.93	0.27	-	75,75,75,75	0
54	MG	BY	103	1/1	0.78	0.22	-	95,95,95,95	0
54	MG	DA	3469	1/1	0.91	0.40	-	83,83,83,83	0
54	MG	CA	1641	1/1	0.71	0.40	-	79,79,79,79	0
54	MG	BB	223	1/1	0.95	0.17	-	94,94,94,94	0
54	MG	BA	3344	1/1	0.86	0.36	-	99,99,99,99	0
54	MG	DA	3235	1/1	0.99	0.20	-	59,59,59,59	0
54	MG	BA	3625	1/1	0.85	0.25	-	53,53,53,53	0
54	MG	BA	3383	1/1	0.90	0.25	-	66,66,66,66	0
54	MG	BA	3199	1/1	0.89	0.33	-	61,61,61,61	0
54	MG	CA	1928	1/1	0.56	0.54	-	117,117,117,117	0
54	MG	AA	1866	1/1	0.65	0.58	-	108,108,108,108	0
54	MG	BA	3179	1/1	0.93	0.21	-	79,79,79,79	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
54	MG	AA	1741	1/1	0.71	0.40	-	112,112,112,112	0
54	MG	BA	3497	1/1	0.92	0.23	-	61,61,61,61	0
54	MG	CA	1802	1/1	0.99	0.03	-	94,94,94,94	0
54	MG	BB	224	1/1	0.83	0.36	-	140,140,140,140	0
54	MG	BA	3247	1/1	0.90	0.25	-	109,109,109,109	0
54	MG	CA	1799	1/1	0.84	0.18	-	109,109,109,109	0
54	MG	DA	3403	1/1	0.85	0.31	-	56,56,56,56	0
54	MG	DA	3579	1/1	0.96	0.54	-	80,80,80,80	0
54	MG	DB	223	1/1	0.83	0.17	-	82,82,82,82	0
54	MG	DA	3068	1/1	0.80	0.53	-	64,64,64,64	0
54	MG	DA	3728	1/1	0.86	0.68	-	83,83,83,83	0
54	MG	BA	3527	1/1	0.79	0.52	-	65,65,65,65	0
54	MG	BA	3138	1/1	0.81	0.31	-	88,88,88,88	0
54	MG	BA	3602	1/1	0.85	0.73	-	98,98,98,98	0
54	MG	BA	3617	1/1	0.90	0.14	-	96,96,96,96	0
54	MG	BA	3275	1/1	0.92	0.24	-	68,68,68,68	0
54	MG	BA	3591	1/1	0.94	0.22	-	99,99,99,99	0
54	MG	DA	3757	1/1	0.97	0.14	-	89,89,89,89	0
54	MG	DA	2910	1/1	0.97	0.68	-	38,38,38,38	0
54	MG	DA	3536	1/1	0.97	0.16	-	92,92,92,92	0
54	MG	DA	3458	1/1	0.86	0.09	-	95,95,95,95	0
54	MG	DL	203	1/1	0.88	0.33	-	66,66,66,66	0
54	MG	BA	3409	1/1	0.97	0.15	-	51,51,51,51	0
54	MG	DA	3819	1/1	0.90	0.58	-	130,130,130,130	0
54	MG	CA	1971	1/1	0.84	0.43	-	86,86,86,86	0
54	MG	DA	3688	1/1	0.78	0.57	-	50,50,50,50	0
54	MG	DA	3440	1/1	0.96	0.46	-	125,125,125,125	0
54	MG	DA	2983	1/1	0.94	0.40	-	52,52,52,52	0
54	MG	CC	303	1/1	0.52	0.23	-	147,147,147,147	0
54	MG	AA	1756	1/1	0.96	0.23	-	111,111,111,111	0
54	MG	DA	2915	1/1	0.92	0.47	-	49,49,49,49	0
54	MG	DA	2905	1/1	0.98	0.33	-	18,18,18,18	0
54	MG	BA	3561	1/1	0.95	0.14	-	75,75,75,75	0
54	MG	BB	215	1/1	0.93	0.39	-	82,82,82,82	0
54	MG	CA	1849	1/1	0.93	0.08	-	167,167,167,167	0
54	MG	BA	2984	1/1	0.90	0.20	-	50,50,50,50	0
54	MG	DA	3163	1/1	0.85	0.31	-	113,113,113,113	0
54	MG	DA	3237	1/1	0.84	0.73	-	64,64,64,64	0
54	MG	BA	3201	1/1	0.82	0.43	-	107,107,107,107	0
54	MG	BA	3346	1/1	0.90	0.15	-	136,136,136,136	0
54	MG	DA	3549	1/1	0.87	0.32	-	115,115,115,115	0
54	MG	DA	3273	1/1	0.90	0.13	-	94,94,94,94	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
54	MG	BA	3386	1/1	0.98	0.36	-	134,134,134,134	0
54	MG	CA	1847	1/1	0.94	0.16	-	58,58,58,58	0
54	MG	DA	3669	1/1	0.72	0.46	-	61,61,61,61	0
54	MG	BA	3091	1/1	0.89	0.22	-	72,72,72,72	0
54	MG	AA	1769	1/1	0.84	0.49	-	118,118,118,118	0
54	MG	BA	3324	1/1	0.76	0.38	-	109,109,109,109	0
54	MG	DA	3592	1/1	0.90	0.12	-	114,114,114,114	0
54	MG	BA	3242	1/1	0.78	0.15	-	112,112,112,112	0
54	MG	DA	3766	1/1	0.64	0.14	-	123,123,123,123	0
54	MG	BA	3299	1/1	0.91	0.15	-	72,72,72,72	0
54	MG	DA	3612	1/1	0.92	0.16	-	74,74,74,74	0
54	MG	DA	3513	1/1	0.99	0.34	-	30,30,30,30	0
54	MG	DA	3229	1/1	0.92	0.22	-	59,59,59,59	0
54	MG	BA	3195	1/1	0.91	0.71	-	91,91,91,91	0
54	MG	DA	3614	1/1	-0.06	0.31	-	128,128,128,128	0
54	MG	DA	3206	1/1	0.92	0.33	-	89,89,89,89	0
54	MG	CA	1712	1/1	0.62	0.25	-	108,108,108,108	0
54	MG	DA	3212	1/1	0.93	0.44	-	75,75,75,75	0
54	MG	CA	1652	1/1	0.96	0.19	-	87,87,87,87	0
54	MG	BA	3593	1/1	0.74	0.14	-	81,81,81,81	0
54	MG	DA	3833	1/1	0.93	0.43	-	73,73,73,73	0
54	MG	CA	1732	1/1	0.70	0.82	-	78,78,78,78	0
54	MG	CA	1945	1/1	0.92	0.64	-	95,95,95,95	0
54	MG	BA	3317	1/1	0.96	0.17	-	84,84,84,84	0
54	MG	DA	3223	1/1	0.79	0.10	-	83,83,83,83	0
54	MG	CA	1663	1/1	0.90	0.31	-	128,128,128,128	0
54	MG	AF	201	1/1	0.84	0.14	-	90,90,90,90	0
54	MG	CA	1743	1/1	0.90	0.14	-	88,88,88,88	0
54	MG	DA	3508	1/1	0.94	0.17	-	70,70,70,70	0
54	MG	CA	1817	1/1	0.95	0.23	-	89,89,89,89	0
54	MG	DA	3687	1/1	0.91	0.18	-	57,57,57,57	0
54	MG	AA	1872	1/1	0.95	0.51	-	88,88,88,88	0
54	MG	CA	1863	1/1	0.81	0.14	-	125,125,125,125	0
54	MG	DA	3675	1/1	0.80	0.62	-	46,46,46,46	0
54	MG	DA	3582	1/1	0.94	0.16	-	85,85,85,85	0
54	MG	DA	3418	1/1	0.74	0.24	-	65,65,65,65	0
54	MG	BB	219	1/1	0.84	0.09	-	69,69,69,69	0
54	MG	DA	3352	1/1	0.95	0.21	-	121,121,121,121	0
54	MG	BA	3401	1/1	0.91	0.27	-	85,85,85,85	0
54	MG	AA	1783	1/1	0.96	0.23	-	123,123,123,123	0
54	MG	BA	3359	1/1	0.97	0.12	-	89,89,89,89	0
54	MG	BA	3502	1/1	0.76	0.12	-	81,81,81,81	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
54	MG	BA	3217	1/1	0.88	0.22	-	55,55,55,55	0
54	MG	CA	1792	1/1	0.86	0.31	-	93,93,93,93	0
54	MG	AA	1648	1/1	0.97	0.49	-	90,90,90,90	0
54	MG	BA	3257	1/1	0.73	1.02	-	108,108,108,108	0
54	MG	BA	3407	1/1	0.88	0.30	-	79,79,79,79	0
54	MG	CE	202	1/1	0.72	0.23	-	80,80,80,80	0
54	MG	DA	3083	1/1	0.95	0.25	-	120,120,120,120	0
54	MG	BA	3263	1/1	0.87	0.32	-	107,107,107,107	0
54	MG	DA	3769	1/1	0.65	0.35	-	90,90,90,90	0
54	MG	BA	3300	1/1	0.96	0.32	-	93,93,93,93	0
54	MG	BA	3285	1/1	0.95	0.11	-	91,91,91,91	0
54	MG	DA	3362	1/1	0.91	0.29	-	78,78,78,78	0
54	MG	DA	3709	1/1	0.93	0.75	-	75,75,75,75	0
54	MG	DA	3814	1/1	0.73	0.17	-	186,186,186,186	0
54	MG	CA	1827	1/1	0.95	0.25	-	82,82,82,82	0
54	MG	BA	3600	1/1	0.79	0.14	-	93,93,93,93	0
54	MG	DA	3405	1/1	0.94	0.87	-	67,67,67,67	0
54	MG	DB	237	1/1	0.74	0.81	-	109,109,109,109	0
54	MG	DA	3565	1/1	0.96	0.17	-	97,97,97,97	0
54	MG	BA	3566	1/1	0.90	0.31	-	72,72,72,72	0
54	MG	CA	1872	1/1	0.93	0.47	-	84,84,84,84	0
54	MG	DA	3644	1/1	0.90	0.50	-	75,75,75,75	0
54	MG	CA	1911	1/1	0.73	0.76	-	90,90,90,90	0
54	MG	BA	3347	1/1	0.89	0.16	-	77,77,77,77	0
54	MG	BA	3005	1/1	0.99	0.76	-	53,53,53,53	0
54	MG	DA	3545	1/1	0.70	0.70	-	94,94,94,94	0
54	MG	BA	3029	1/1	0.97	0.33	-	41,41,41,41	0
54	MG	BA	3587	1/1	0.90	0.27	-	77,77,77,77	0
54	MG	BA	3093	1/1	0.89	0.39	-	67,67,67,67	0
54	MG	DA	3350	1/1	0.92	0.09	-	95,95,95,95	0
54	MG	DA	2939	1/1	0.98	0.40	-	21,21,21,21	0
54	MG	BA	3598	1/1	0.91	0.10	-	100,100,100,100	0
54	MG	BA	3077	1/1	0.82	0.51	-	55,55,55,55	0
54	MG	DA	3804	1/1	0.93	0.32	-	72,72,72,72	0
54	MG	BA	2952	1/1	0.93	0.19	-	59,59,59,59	0
54	MG	DA	3245	1/1	0.92	0.79	-	100,100,100,100	0
54	MG	AA	1686	1/1	0.92	0.47	-	65,65,65,65	0
54	MG	BA	3130	1/1	0.93	0.21	-	47,47,47,47	0
54	MG	DA	3093	1/1	0.88	0.35	-	47,47,47,47	0
54	MG	AA	1766	1/1	0.89	0.24	-	106,106,106,106	0
54	MG	BA	3108	1/1	0.81	0.50	-	72,72,72,72	0
54	MG	CA	1643	1/1	0.77	0.19	-	70,70,70,70	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
54	MG	AA	1859	1/1	0.89	0.21	-	87,87,87,87	0
54	MG	BB	213	1/1	0.90	0.41	-	111,111,111,111	0
54	MG	BA	3453	1/1	0.96	0.27	-	27,27,27,27	0
54	MG	DA	3264	1/1	0.89	0.12	-	78,78,78,78	0
54	MG	DA	3406	1/1	0.93	0.20	-	59,59,59,59	0
54	MG	DA	3115	1/1	0.96	0.18	-	58,58,58,58	0
54	MG	BA	2944	1/1	0.92	0.51	-	47,47,47,47	0
54	MG	CA	1938	1/1	0.80	0.20	-	81,81,81,81	0
54	MG	AV	6202	1/1	0.71	0.08	-	97,97,97,97	0
54	MG	DA	3230	1/1	0.97	0.25	-	80,80,80,80	0
54	MG	DA	2903	1/1	0.99	0.24	-	16,16,16,16	0
54	MG	CA	1605	1/1	0.96	0.64	-	54,54,54,54	0
54	MG	BA	3094	1/1	0.76	0.27	-	63,63,63,63	0
54	MG	DA	3556	1/1	0.96	0.15	-	110,110,110,110	0
54	MG	BA	3113	1/1	0.97	0.53	-	59,59,59,59	0
54	MG	DA	2937	1/1	0.94	0.43	-	43,43,43,43	0
54	MG	DA	3553	1/1	0.94	0.14	-	91,91,91,91	0
54	MG	DA	3084	1/1	0.91	0.85	-	65,65,65,65	0
54	MG	BA	3444	1/1	0.97	0.32	-	26,26,26,26	0
54	MG	DA	3138	1/1	0.93	0.30	-	81,81,81,81	0
54	MG	BA	3111	1/1	0.91	0.72	-	64,64,64,64	0
54	MG	DA	3397	1/1	0.81	0.92	-	124,124,124,124	0
54	MG	DD	304	1/1	0.73	0.96	-	94,94,94,94	0
54	MG	BA	3537	1/1	0.88	0.19	-	60,60,60,60	0
54	MG	CA	1831	1/1	0.93	0.44	-	90,90,90,90	0
54	MG	AD	304	1/1	0.68	0.44	-	113,113,113,113	0
54	MG	DA	3512	1/1	0.91	0.35	-	159,159,159,159	0
54	MG	DA	3711	1/1	0.83	0.44	-	84,84,84,84	0
54	MG	CA	1721	1/1	0.88	0.91	-	118,118,118,118	0
54	MG	BA	3255	1/1	0.86	0.57	-	101,101,101,101	0
54	MG	BA	3063	1/1	0.96	0.92	-	85,85,85,85	0
54	MG	BA	3633	1/1	0.63	0.49	-	88,88,88,88	0
54	MG	DA	3085	1/1	0.91	0.46	-	68,68,68,68	0
54	MG	DA	2956	1/1	0.90	0.43	-	44,44,44,44	0
54	MG	BA	3182	1/1	0.95	0.19	-	78,78,78,78	0
54	MG	AA	1676	1/1	0.95	0.14	-	90,90,90,90	0
54	MG	DA	3122	1/1	0.64	0.27	-	93,93,93,93	0
54	MG	DA	3054	1/1	0.95	1.00	-	72,72,72,72	0
54	MG	BA	3590	1/1	0.65	0.34	-	119,119,119,119	0
54	MG	DA	3813	1/1	0.94	0.15	-	50,50,50,50	0
54	MG	BA	3081	1/1	0.91	0.21	-	67,67,67,67	0
54	MG	AA	1689	1/1	0.92	0.62	-	75,75,75,75	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
54	MG	CA	1957	1/1	0.97	0.43	-	73,73,73,73	0
54	MG	AA	1823	1/1	0.79	0.63	-	148,148,148,148	0
54	MG	DA	3239	1/1	0.93	0.18	-	126,126,126,126	0
54	MG	AL	202	1/1	0.84	0.12	-	67,67,67,67	0
54	MG	AA	1770	1/1	0.91	0.21	-	76,76,76,76	0
54	MG	AA	1762	1/1	0.90	0.55	-	91,91,91,91	0
54	MG	AA	1797	1/1	0.89	0.48	-	72,72,72,72	0
54	MG	AA	1768	1/1	0.95	0.80	-	87,87,87,87	0
54	MG	DA	3699	1/1	0.58	0.12	-	96,96,96,96	0
54	MG	AA	1771	1/1	0.71	1.16	-	111,111,111,111	0
54	MG	DA	3542	1/1	0.95	0.24	-	65,65,65,65	0
54	MG	BA	3627	1/1	0.82	0.34	-	67,67,67,67	0
54	MG	DA	3659	1/1	0.87	1.14	-	61,61,61,61	0
54	MG	BA	2978	1/1	0.76	0.71	-	58,58,58,58	0
54	MG	CA	1907	1/1	0.88	0.44	-	87,87,87,87	0
54	MG	CA	1897	1/1	0.76	0.81	-	74,74,74,74	0
54	MG	DA	3188	1/1	0.81	0.53	-	69,69,69,69	0
54	MG	BA	3224	1/1	0.97	0.25	-	97,97,97,97	0
54	MG	BA	3543	1/1	0.86	0.30	-	76,76,76,76	0
54	MG	BA	3032	1/1	0.82	0.37	-	76,76,76,76	0
54	MG	DA	3200	1/1	0.95	0.30	-	97,97,97,97	0
54	MG	CA	1820	1/1	0.73	1.15	-	99,99,99,99	0
54	MG	BA	3416	1/1	0.77	0.12	-	83,83,83,83	0
54	MG	DK	203	1/1	0.83	0.24	-	73,73,73,73	0
54	MG	DA	3236	1/1	0.96	0.15	-	75,75,75,75	0
54	MG	BA	3296	1/1	0.97	0.22	-	71,71,71,71	0
54	MG	DA	2967	1/1	0.96	0.33	-	34,34,34,34	0
54	MG	CA	1836	1/1	0.98	0.43	-	98,98,98,98	0
54	MG	AA	1692	1/1	0.93	0.16	-	89,89,89,89	0
54	MG	DA	3821	1/1	0.92	0.34	-	79,79,79,79	0
54	MG	DA	3732	1/1	0.82	0.32	-	81,81,81,81	0
54	MG	BA	3559	1/1	0.79	0.76	-	70,70,70,70	0
54	MG	DA	3253	1/1	0.90	0.22	-	70,70,70,70	0
54	MG	CA	1776	1/1	0.90	0.13	-	118,118,118,118	0
54	MG	DD	302	1/1	0.89	0.21	-	61,61,61,61	0
54	MG	DA	3839	1/1	0.68	0.30	-	86,86,86,86	0
54	MG	BA	3279	1/1	0.99	0.06	-	85,85,85,85	0
54	MG	CA	1841	1/1	0.90	0.18	-	95,95,95,95	0
54	MG	AA	1671	1/1	0.95	0.13	-	74,74,74,74	0
54	MG	DB	208	1/1	0.92	0.16	-	86,86,86,86	0
54	MG	BA	3601	1/1	0.96	0.20	-	76,76,76,76	0
54	MG	CA	1934	1/1	0.95	0.12	-	80,80,80,80	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
54	MG	DA	2946	1/1	0.97	0.51	-	25,25,25,25	0
54	MG	AA	1816	1/1	0.88	0.88	-	89,89,89,89	0
54	MG	AA	1799	1/1	0.81	0.52	-	58,58,58,58	0
54	MG	CA	1734	1/1	0.74	0.87	-	90,90,90,90	0
54	MG	AA	1750	1/1	0.91	0.30	-	128,128,128,128	0
54	MG	AA	1818	1/1	0.75	0.23	-	76,76,76,76	0
54	MG	CA	1757	1/1	0.97	0.18	-	102,102,102,102	0
54	MG	DA	3183	1/1	0.98	0.33	-	76,76,76,76	0
54	MG	DA	3611	1/1	0.95	0.77	-	74,74,74,74	0
54	MG	BA	3282	1/1	0.76	0.40	-	98,98,98,98	0
54	MG	DA	3358	1/1	0.96	0.30	-	83,83,83,83	0
54	MG	BA	3583	1/1	0.96	0.70	-	105,105,105,105	0
54	MG	BA	3354	1/1	0.53	0.56	-	123,123,123,123	0
54	MG	DA	3049	1/1	0.98	0.52	-	43,43,43,43	0
54	MG	BA	3314	1/1	0.96	0.20	-	69,69,69,69	0
54	MG	CA	1959	1/1	0.96	0.09	-	106,106,106,106	0
54	MG	BA	3665	1/1	0.91	0.25	-	123,123,123,123	0
54	MG	BA	3516	1/1	0.66	0.39	-	100,100,100,100	0
54	MG	DA	3824	1/1	0.94	0.25	-	113,113,113,113	0
54	MG	AA	1848	1/1	0.88	0.34	-	77,77,77,77	0
54	MG	DA	3523	1/1	0.96	0.13	-	77,77,77,77	0
54	MG	BB	210	1/1	0.91	0.29	-	82,82,82,82	0
54	MG	DA	3348	1/1	0.87	0.10	-	83,83,83,83	0
54	MG	CA	1925	1/1	0.86	0.37	-	77,77,77,77	0
54	MG	DA	3786	1/1	0.79	0.90	-	104,104,104,104	0
54	MG	DA	3425	1/1	0.80	0.25	-	113,113,113,113	0
54	MG	DA	3474	1/1	0.62	1.06	-	81,81,81,81	0
54	MG	DA	3282	1/1	0.90	0.55	-	69,69,69,69	0
54	MG	CB	301	1/1	0.82	0.22	-	115,115,115,115	0
54	MG	AI	201	1/1	0.92	0.19	-	70,70,70,70	0
54	MG	CA	1699	1/1	0.97	0.10	-	76,76,76,76	0
54	MG	AA	1702	1/1	0.95	0.15	-	56,56,56,56	0
54	MG	CA	1868	1/1	0.94	0.30	-	71,71,71,71	0
54	MG	BA	3610	1/1	0.88	0.90	-	80,80,80,80	0
54	MG	AA	1754	1/1	0.95	0.16	-	98,98,98,98	0
54	MG	CA	1661	1/1	0.94	0.11	-	80,80,80,80	0
54	MG	CA	1830	1/1	0.80	0.26	-	82,82,82,82	0
54	MG	AA	1725	1/1	0.94	0.24	-	93,93,93,93	0
54	MG	CA	1794	1/1	0.99	0.12	-	122,122,122,122	0
54	MG	CA	1825	1/1	0.92	0.46	-	97,97,97,97	0
54	MG	DA	2944	1/1	0.93	0.43	-	25,25,25,25	0
54	MG	DA	3487	1/1	0.92	0.40	-	66,66,66,66	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
54	MG	BA	3293	1/1	0.96	0.24	-	60,60,60,60	0
54	MG	BA	3135	1/1	0.83	0.24	-	61,61,61,61	0
54	MG	DA	3216	1/1	0.92	0.54	-	91,91,91,91	0
54	MG	CO	104	1/1	0.63	0.24	-	99,99,99,99	0
54	MG	AA	1633	1/1	0.83	0.29	-	123,123,123,123	0
54	MG	AA	1851	1/1	0.91	0.62	-	93,93,93,93	0
54	MG	DA	3712	1/1	0.82	0.20	-	41,41,41,41	0
54	MG	BA	3538	1/1	0.93	0.25	-	75,75,75,75	0
54	MG	BA	2987	1/1	0.85	0.06	-	79,79,79,79	0
54	MG	DA	3147	1/1	0.88	0.47	-	86,86,86,86	0
54	MG	BC	301	1/1	0.86	0.80	-	60,60,60,60	0
54	MG	BA	3638	1/1	0.89	0.85	-	127,127,127,127	0
54	MG	BA	3398	1/1	0.93	0.28	-	74,74,74,74	0
54	MG	DA	3586	1/1	0.89	0.32	-	79,79,79,79	0
54	MG	CA	1874	1/1	0.78	0.38	-	76,76,76,76	0
54	MG	DA	3031	1/1	0.90	0.62	-	79,79,79,79	0
54	MG	AA	1753	1/1	0.85	0.60	-	103,103,103,103	0
54	MG	BA	3229	1/1	0.75	0.27	-	90,90,90,90	0
54	MG	DB	221	1/1	0.95	0.29	-	98,98,98,98	0
54	MG	AA	1687	1/1	0.85	0.70	-	108,108,108,108	0
54	MG	BA	3312	1/1	0.97	0.69	-	71,71,71,71	0
54	MG	DA	3783	1/1	0.96	0.13	-	63,63,63,63	0
54	MG	BA	3392	1/1	0.80	0.27	-	108,108,108,108	0
54	MG	DA	3430	1/1	0.96	0.27	-	103,103,103,103	0
54	MG	CD	302	1/1	0.83	0.13	-	74,74,74,74	0
54	MG	CA	1774	1/1	0.90	0.27	-	79,79,79,79	0
54	MG	AA	1791	1/1	0.92	0.08	-	106,106,106,106	0
54	MG	DA	3208	1/1	0.97	0.28	-	56,56,56,56	0
54	MG	BA	3302	1/1	0.93	0.34	-	84,84,84,84	0
54	MG	CA	1913	1/1	0.89	0.19	-	109,109,109,109	0
54	MG	D1	101	1/1	0.65	0.39	-	80,80,80,80	0
54	MG	BA	3609	1/1	0.93	0.16	-	65,65,65,65	0
54	MG	CA	1954	1/1	0.85	0.08	-	89,89,89,89	0
54	MG	BA	3552	1/1	0.87	0.11	-	99,99,99,99	0
54	MG	CA	1812	1/1	0.93	0.74	-	92,92,92,92	0
54	MG	DA	3578	1/1	0.69	0.29	-	96,96,96,96	0
54	MG	BA	3372	1/1	0.74	1.46	-	123,123,123,123	0
54	MG	BB	206	1/1	0.83	0.47	-	125,125,125,125	0
54	MG	BA	3389	1/1	0.92	0.30	-	66,66,66,66	0
54	MG	DA	3530	1/1	0.88	0.32	-	71,71,71,71	0
54	MG	AA	1796	1/1	0.91	0.62	-	106,106,106,106	0
54	MG	BA	3680	1/1	0.71	0.76	-	97,97,97,97	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
54	MG	DA	3518	1/1	0.96	0.24	-	102,102,102,102	0
54	MG	DA	3155	1/1	0.78	0.29	-	70,70,70,70	0
54	MG	DA	3370	1/1	0.86	1.28	-	92,92,92,92	0
54	MG	BA	3298	1/1	0.97	0.22	-	126,126,126,126	0
54	MG	DA	3808	1/1	0.85	0.32	-	116,116,116,116	0
54	MG	DA	3744	1/1	0.92	0.24	-	64,64,64,64	0
54	MG	DA	3714	1/1	0.90	0.07	-	96,96,96,96	0
54	MG	BA	3200	1/1	0.98	0.26	-	93,93,93,93	0
54	MG	BA	3465	1/1	0.91	0.32	-	31,31,31,31	0
54	MG	AA	1728	1/1	0.80	0.29	-	71,71,71,71	0
54	MG	BA	2969	1/1	0.94	0.56	-	69,69,69,69	0
54	MG	BA	3284	1/1	0.76	0.24	-	95,95,95,95	0
54	MG	BA	3212	1/1	0.89	0.27	-	67,67,67,67	0
54	MG	BA	3318	1/1	0.87	0.48	-	73,73,73,73	0
54	MG	DA	3278	1/1	0.74	0.74	-	99,99,99,99	0
54	MG	CA	1604	1/1	0.96	0.48	-	43,43,43,43	0
54	MG	BA	3668	1/1	0.68	0.47	-	97,97,97,97	0
54	MG	DA	3494	1/1	0.88	0.29	-	63,63,63,63	0
54	MG	DA	3488	1/1	0.96	0.29	-	102,102,102,102	0
54	MG	BA	3436	1/1	0.97	0.53	-	38,38,38,38	0
54	MG	DA	3433	1/1	0.93	0.25	-	67,67,67,67	0
54	MG	CL	202	1/1	0.97	0.07	-	105,105,105,105	0
54	MG	BA	3115	1/1	0.87	0.44	-	76,76,76,76	0
54	MG	DA	3324	1/1	0.92	0.25	-	95,95,95,95	0
54	MG	AA	1792	1/1	0.79	0.10	-	101,101,101,101	0
54	MG	DA	3075	1/1	0.84	0.19	-	50,50,50,50	0
54	MG	DA	3661	1/1	0.95	0.64	-	98,98,98,98	0
54	MG	CA	1636	1/1	0.95	0.08	-	98,98,98,98	0
54	MG	DA	3211	1/1	0.75	0.37	-	64,64,64,64	0
54	MG	BA	3657	1/1	0.82	0.43	-	84,84,84,84	0
54	MG	BA	3155	1/1	0.75	0.93	-	99,99,99,99	0
54	MG	DA	3325	1/1	0.92	0.25	-	66,66,66,66	0
54	MG	BA	3420	1/1	0.72	0.53	-	94,94,94,94	0
54	MG	DA	3076	1/1	0.94	0.13	-	78,78,78,78	0
54	MG	CA	1808	1/1	0.94	0.14	-	98,98,98,98	0
54	MG	CA	1864	1/1	0.95	0.21	-	90,90,90,90	0
54	MG	BA	3422	1/1	0.91	0.18	-	93,93,93,93	0
54	MG	CA	1664	1/1	0.98	0.09	-	85,85,85,85	0
54	MG	DA	2993	1/1	0.86	0.38	-	47,47,47,47	0
54	MG	BA	3395	1/1	0.94	0.40	-	84,84,84,84	0
54	MG	AA	1654	1/1	0.94	0.28	-	99,99,99,99	0
54	MG	BA	3577	1/1	0.95	0.32	-	142,142,142,142	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
54	MG	AA	1706	1/1	0.96	0.46	-	78,78,78,78	0
54	MG	DB	240	1/1	0.90	0.17	-	123,123,123,123	0
54	MG	AA	1742	1/1	0.93	0.14	-	90,90,90,90	0
54	MG	BA	2986	1/1	0.96	0.72	-	54,54,54,54	0
54	MG	DA	3302	1/1	0.87	0.73	-	102,102,102,102	0
54	MG	DA	3380	1/1	0.80	0.21	-	84,84,84,84	0
54	MG	BA	3181	1/1	0.92	0.21	-	62,62,62,62	0
54	MG	DA	3393	1/1	0.88	0.40	-	77,77,77,77	0
54	MG	BA	3132	1/1	0.68	0.26	-	66,66,66,66	0
54	MG	AV	6208	1/1	0.96	0.38	-	79,79,79,79	0
54	MG	CE	203	1/1	0.70	0.12	-	79,79,79,79	0
54	MG	CG	203	1/1	0.80	0.22	-	100,100,100,100	0
54	MG	BA	3028	1/1	0.85	1.61	-	61,61,61,61	0
54	MG	DA	2912	1/1	0.94	0.41	-	19,19,19,19	0
54	MG	BR	201	1/1	0.89	0.07	-	75,75,75,75	0
54	MG	AA	1677	1/1	0.97	0.17	-	50,50,50,50	0
54	MG	DA	3723	1/1	0.88	0.31	-	67,67,67,67	0
54	MG	DA	3682	1/1	0.91	0.23	-	59,59,59,59	0
54	MG	AA	1855	1/1	0.93	0.34	-	81,81,81,81	0
54	MG	DA	3201	1/1	0.98	0.05	-	46,46,46,46	0
54	MG	BA	3355	1/1	0.96	0.22	-	72,72,72,72	0
54	MG	BA	3231	1/1	0.94	0.29	-	64,64,64,64	0
54	MG	CB	302	1/1	0.92	0.21	-	112,112,112,112	0
54	MG	CA	1892	1/1	0.90	0.65	-	114,114,114,114	0
54	MG	BA	3283	1/1	0.84	0.26	-	100,100,100,100	0
54	MG	BA	3539	1/1	0.85	0.42	-	97,97,97,97	0
54	MG	BA	3345	1/1	0.92	0.26	-	71,71,71,71	0
54	MG	DA	3148	1/1	0.92	0.12	-	94,94,94,94	0
54	MG	CA	1722	1/1	0.74	0.30	-	116,116,116,116	0
54	MG	BA	3331	1/1	0.95	0.86	-	66,66,66,66	0
54	MG	AA	1878	1/1	0.77	0.09	-	92,92,92,92	0
54	MG	DA	3456	1/1	0.96	0.47	-	121,121,121,121	0
54	MG	CA	1923	1/1	0.93	0.31	-	74,74,74,74	0
54	MG	AA	1623	1/1	0.66	0.16	-	99,99,99,99	0
54	MG	DA	3845	1/1	0.90	0.23	-	94,94,94,94	0
54	MG	BA	3565	1/1	0.97	0.25	-	104,104,104,104	0
54	MG	DA	3269	1/1	0.93	0.19	-	109,109,109,109	0
54	MG	DA	3219	1/1	0.90	0.14	-	60,60,60,60	0
54	MG	BA	3511	1/1	0.82	1.01	-	78,78,78,78	0
54	MG	AA	1819	1/1	0.81	0.45	-	73,73,73,73	0
54	MG	AA	1801	1/1	0.93	0.89	-	71,71,71,71	0
54	MG	AA	1747	1/1	0.88	0.45	-	116,116,116,116	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
54	MG	DA	3716	1/1	0.82	0.58	-	78,78,78,78	0
54	MG	BA	2925	1/1	0.98	0.29	-	12,12,12,12	0
54	MG	DA	3368	1/1	0.88	0.25	-	98,98,98,98	0
54	MG	DB	202	1/1	0.96	0.30	-	33,33,33,33	0
54	MG	DA	3794	1/1	0.92	0.23	-	104,104,104,104	0
54	MG	DA	3809	1/1	0.76	1.09	-	84,84,84,84	0
54	MG	CA	1706	1/1	0.89	0.82	-	106,106,106,106	0
54	MG	BA	2953	1/1	0.96	0.18	-	45,45,45,45	0
54	MG	DA	3411	1/1	0.79	0.51	-	68,68,68,68	0
54	MG	DA	3290	1/1	0.90	0.15	-	61,61,61,61	0
54	MG	AA	1803	1/1	0.83	0.49	-	69,69,69,69	0
54	MG	CA	1895	1/1	0.92	0.73	-	88,88,88,88	0
54	MG	AA	1789	1/1	0.82	0.41	-	106,106,106,106	0
54	MG	BA	3207	1/1	0.94	0.37	-	70,70,70,70	0
54	MG	DA	3547	1/1	0.98	0.40	-	115,115,115,115	0
54	MG	DA	3112	1/1	0.94	0.27	-	91,91,91,91	0
54	MG	AA	1765	1/1	0.94	0.14	-	91,91,91,91	0
54	MG	DA	3191	1/1	0.89	0.23	-	62,62,62,62	0
54	MG	DA	3651	1/1	0.91	0.37	-	51,51,51,51	0
54	MG	AA	1794	1/1	0.94	0.15	-	112,112,112,112	0
54	MG	DA	3507	1/1	0.97	0.65	-	70,70,70,70	0
54	MG	BA	2959	1/1	0.96	0.31	-	40,40,40,40	0
54	MG	DA	3576	1/1	0.94	0.49	-	99,99,99,99	0
54	MG	BA	3292	1/1	0.88	0.24	-	71,71,71,71	0
54	MG	BA	3335	1/1	0.85	0.89	-	80,80,80,80	0
54	MG	DA	3081	1/1	0.83	0.68	-	67,67,67,67	0
54	MG	DB	218	1/1	0.94	0.10	-	113,113,113,113	0
54	MG	DA	3798	1/1	0.95	0.43	-	88,88,88,88	0
54	MG	DA	3315	1/1	0.94	0.40	-	58,58,58,58	0
54	MG	DA	3450	1/1	0.76	0.72	-	81,81,81,81	0
54	MG	DA	3642	1/1	0.91	0.45	-	52,52,52,52	0
54	MG	CA	1834	1/1	0.94	0.17	-	81,81,81,81	0
54	MG	DA	3145	1/1	0.94	0.16	-	100,100,100,100	0
54	MG	CA	1900	1/1	0.81	0.23	-	86,86,86,86	0
54	MG	DA	3133	1/1	0.98	0.18	-	114,114,114,114	0
54	MG	DF	201	1/1	0.79	0.32	-	70,70,70,70	0
54	MG	CA	1844	1/1	0.85	0.41	-	78,78,78,78	0
54	MG	DA	3822	1/1	0.91	0.11	-	70,70,70,70	0
54	MG	AA	1602	1/1	0.95	0.89	-	45,45,45,45	0
54	MG	AA	1864	1/1	0.83	0.17	-	75,75,75,75	0
54	MG	BA	3209	1/1	0.92	0.60	-	81,81,81,81	0
54	MG	BA	2951	1/1	0.90	0.84	-	53,53,53,53	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
54	MG	BA	3168	1/1	0.98	0.27	-	20,20,20,20	0
54	MG	DB	222	1/1	0.95	0.19	-	137,137,137,137	0
54	MG	BA	3472	1/1	0.95	0.51	-	65,65,65,65	0
54	MG	DA	3187	1/1	0.92	0.23	-	114,114,114,114	0
54	MG	AA	1703	1/1	0.93	0.45	-	63,63,63,63	0
54	MG	BA	3614	1/1	0.68	0.11	-	112,112,112,112	0
54	MG	DA	3841	1/1	0.81	0.40	-	80,80,80,80	0
54	MG	DA	3587	1/1	0.74	0.12	-	113,113,113,113	0
54	MG	DA	3313	1/1	0.89	0.54	-	78,78,78,78	0
54	MG	CA	1779	1/1	0.96	0.28	-	79,79,79,79	0
54	MG	BA	3521	1/1	0.89	0.19	-	88,88,88,88	0
54	MG	CA	1611	1/1	0.94	0.30	-	128,128,128,128	0
54	MG	BA	3528	1/1	0.87	0.22	-	73,73,73,73	0
54	MG	BY	101	1/1	0.73	0.34	-	42,42,42,42	0
54	MG	DA	3510	1/1	0.96	0.23	-	64,64,64,64	0
54	MG	DA	3213	1/1	0.99	0.25	-	59,59,59,59	0
54	MG	BB	214	1/1	0.88	0.65	-	84,84,84,84	0
54	MG	AA	1712	1/1	0.66	0.19	-	105,105,105,105	0
54	MG	AA	1716	1/1	0.80	0.32	-	96,96,96,96	0
54	MG	DA	2914	1/1	0.97	0.44	-	23,23,23,23	0
54	MG	CA	1618	1/1	0.91	0.69	-	84,84,84,84	0
54	MG	DA	3759	1/1	0.79	0.53	-	97,97,97,97	0
54	MG	BT	101	1/1	0.86	0.89	-	101,101,101,101	0
54	MG	CV	6201	1/1	0.72	0.10	-	95,95,95,95	0
54	MG	CA	1855	1/1	0.93	0.31	-	107,107,107,107	0
54	MG	CA	1625	1/1	0.86	0.51	-	67,67,67,67	0
54	MG	DA	3670	1/1	0.85	0.61	-	83,83,83,83	0
54	MG	DA	3134	1/1	0.92	0.75	-	69,69,69,69	0
54	MG	BA	3548	1/1	0.83	0.32	-	61,61,61,61	0
54	MG	CA	1851	1/1	0.95	0.48	-	103,103,103,103	0
54	MG	AA	1655	1/1	0.90	0.73	-	58,58,58,58	0
54	MG	DA	3708	1/1	0.80	0.18	-	63,63,63,63	0
54	MG	BA	3160	1/1	0.93	0.63	-	87,87,87,87	0
54	MG	DA	3596	1/1	0.82	0.81	-	113,113,113,113	0
54	MG	CA	1914	1/1	0.89	0.13	-	97,97,97,97	0
54	MG	DA	2964	1/1	0.95	0.46	-	45,45,45,45	0
54	MG	DA	3420	1/1	0.82	0.21	-	110,110,110,110	0
54	MG	DA	3136	1/1	0.97	0.21	-	57,57,57,57	0
54	MG	DA	3697	1/1	0.75	0.34	-	92,92,92,92	0
54	MG	DA	2981	1/1	0.95	0.30	-	40,40,40,40	0
54	MG	DA	3857	1/1	0.91	0.23	-	98,98,98,98	0
54	MG	DA	3396	1/1	0.96	0.28	-	78,78,78,78	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
54	MG	BA	3603	1/1	0.91	0.22	-	86,86,86,86	0
54	MG	DA	3169	1/1	0.94	0.30	-	54,54,54,54	0
54	MG	BA	3531	1/1	0.68	0.14	-	91,91,91,91	0
54	MG	DA	3855	1/1	0.86	0.33	-	104,104,104,104	0
54	MG	BA	2962	1/1	0.77	0.30	-	54,54,54,54	0
54	MG	DA	3571	1/1	0.81	1.03	-	72,72,72,72	0
54	MG	CA	1687	1/1	0.91	1.14	-	63,63,63,63	0
54	MG	DA	3404	1/1	0.90	0.28	-	79,79,79,79	0
54	MG	DA	3285	1/1	0.93	0.52	-	65,65,65,65	0
54	MG	BA	3184	1/1	0.79	0.61	-	67,67,67,67	0
54	MG	DA	3539	1/1	0.62	0.16	-	81,81,81,81	0
54	MG	DA	3829	1/1	0.88	0.08	-	96,96,96,96	0
54	MG	AA	1621	1/1	0.74	0.66	-	71,71,71,71	0
54	MG	BB	201	1/1	0.94	0.48	-	63,63,63,63	0
54	MG	DV	201	1/1	0.84	0.20	-	90,90,90,90	0
54	MG	DA	3305	1/1	0.96	0.28	-	99,99,99,99	0
54	MG	BA	3397	1/1	0.52	0.40	-	125,125,125,125	0
54	MG	BA	3554	1/1	0.93	0.94	-	90,90,90,90	0
54	MG	AA	1617	1/1	0.83	0.59	-	84,84,84,84	0
54	MG	BA	3514	1/1	0.85	0.19	-	83,83,83,83	0
54	MG	AT	201	1/1	0.97	0.09	-	103,103,103,103	0
54	MG	DA	3078	1/1	0.97	0.44	-	48,48,48,48	0
54	MG	BA	3271	1/1	0.95	0.19	-	83,83,83,83	0
54	MG	BA	3445	1/1	0.84	0.46	-	61,61,61,61	0
54	MG	AA	1782	1/1	0.92	0.25	-	104,104,104,104	0
54	MG	DA	3551	1/1	0.91	0.22	-	80,80,80,80	0
54	MG	CA	1947	1/1	0.85	0.21	-	136,136,136,136	0
54	MG	BA	3635	1/1	0.90	0.28	-	113,113,113,113	0
54	MG	BA	3326	1/1	0.96	0.20	-	82,82,82,82	0
54	MG	AA	1865	1/1	0.88	0.12	-	107,107,107,107	0
54	MG	AA	1808	1/1	0.52	0.21	-	97,97,97,97	0
54	MG	DA	3727	1/1	0.85	0.33	-	88,88,88,88	0
54	MG	DA	3760	1/1	0.92	0.27	-	97,97,97,97	0
54	MG	AA	1629	1/1	0.83	0.43	-	66,66,66,66	0
54	MG	AA	1735	1/1	0.85	0.06	-	68,68,68,68	0
54	MG	CA	1735	1/1	0.62	0.52	-	77,77,77,77	0
54	MG	DB	211	1/1	0.96	0.07	-	65,65,65,65	0
54	MG	CC	304	1/1	0.83	0.21	-	121,121,121,121	0
54	MG	BA	2998	1/1	0.95	0.10	-	75,75,75,75	0
54	MG	DA	3203	1/1	0.92	0.10	-	67,67,67,67	0
54	MG	CA	1654	1/1	0.89	0.20	-	78,78,78,78	0
54	MG	DA	3060	1/1	0.96	0.48	-	77,77,77,77	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
54	MG	DA	3328	1/1	0.93	0.33	-	85,85,85,85	0
54	MG	BA	3007	1/1	0.95	0.06	-	98,98,98,98	0
54	MG	BA	3036	1/1	0.91	1.45	-	57,57,57,57	0
54	MG	DA	3751	1/1	0.93	0.21	-	76,76,76,76	0
54	MG	BA	3586	1/1	0.85	0.28	-	97,97,97,97	0
54	MG	AA	1751	1/1	0.88	0.13	-	88,88,88,88	0
54	MG	AA	1675	1/1	0.93	0.27	-	82,82,82,82	0
54	MG	BA	3675	1/1	0.92	0.09	-	120,120,120,120	0
54	MG	CA	1724	1/1	0.48	0.58	-	124,124,124,124	0
54	MG	DA	3643	1/1	0.94	0.48	-	40,40,40,40	0
54	MG	DA	3438	1/1	0.94	0.22	-	54,54,54,54	0
54	MG	DA	3538	1/1	0.73	0.70	-	66,66,66,66	0
54	MG	BA	3185	1/1	0.80	0.40	-	85,85,85,85	0
54	MG	BA	3661	1/1	0.92	0.09	-	82,82,82,82	0
54	MG	BA	3240	1/1	0.93	0.38	-	157,157,157,157	0
54	MG	DA	3058	1/1	0.97	0.14	-	56,56,56,56	0
54	MG	DA	3445	1/1	0.88	0.15	-	137,137,137,137	0
54	MG	CA	1882	1/1	0.47	0.14	-	135,135,135,135	0
54	MG	CA	1929	1/1	0.87	0.48	-	95,95,95,95	0
54	MG	BA	3488	1/1	0.92	0.36	-	53,53,53,53	0
54	MG	AA	1763	1/1	0.92	0.06	-	121,121,121,121	0
54	MG	BA	3082	1/1	0.95	0.14	-	38,38,38,38	0
54	MG	BB	211	1/1	0.83	0.45	-	85,85,85,85	0
54	MG	AA	1608	1/1	0.94	0.41	-	46,46,46,46	0
54	MG	CA	1621	1/1	0.91	0.17	-	67,67,67,67	0
54	MG	CA	1936	1/1	0.94	0.40	-	102,102,102,102	0
54	MG	CA	1698	1/1	0.91	0.16	-	123,123,123,123	0
54	MG	DA	3701	1/1	0.91	0.21	-	56,56,56,56	0
54	MG	CA	1614	1/1	0.76	0.55	-	66,66,66,66	0
54	MG	BA	3644	1/1	0.90	0.37	-	86,86,86,86	0
54	MG	BA	3594	1/1	0.62	0.17	-	129,129,129,129	0
54	MG	DA	3222	1/1	0.88	0.49	-	112,112,112,112	0
54	MG	AA	1646	1/1	0.92	0.67	-	89,89,89,89	0
54	MG	DA	3398	1/1	0.98	0.17	-	80,80,80,80	0
54	MG	CA	1843	1/1	0.78	0.16	-	109,109,109,109	0
54	MG	DA	2917	1/1	0.81	0.18	-	71,71,71,71	0
54	MG	DA	3095	1/1	0.89	0.55	-	72,72,72,72	0
54	MG	BM	201	1/1	0.89	0.28	-	50,50,50,50	0
54	MG	CA	1958	1/1	0.65	0.62	-	125,125,125,125	0
54	MG	BA	3213	1/1	0.91	0.15	-	102,102,102,102	0
54	MG	BA	3315	1/1	0.87	0.15	-	84,84,84,84	0
54	MG	DA	3773	1/1	0.91	0.50	-	37,37,37,37	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
54	MG	CA	1645	1/1	0.96	0.15	-	80,80,80,80	0
54	MG	DA	3385	1/1	0.75	0.46	-	69,69,69,69	0
54	MG	CA	1876	1/1	0.91	0.17	-	69,69,69,69	0
54	MG	DA	2969	1/1	0.95	0.50	-	39,39,39,39	0
54	MG	CA	1685	1/1	0.56	0.33	-	109,109,109,109	0
54	MG	BA	3286	1/1	0.89	0.20	-	68,68,68,68	0
54	MG	BA	3501	1/1	0.81	1.80	-	105,105,105,105	0
54	MG	DA	2977	1/1	0.98	0.57	-	49,49,49,49	0
54	MG	DA	3626	1/1	0.98	0.21	-	106,106,106,106	0
54	MG	BC	302	1/1	0.67	0.48	-	82,82,82,82	0
54	MG	DA	3555	1/1	0.95	0.54	-	61,61,61,61	0
54	MG	DA	3446	1/1	0.96	0.21	-	57,57,57,57	0
54	MG	AA	1781	1/1	0.87	0.20	-	92,92,92,92	0
54	MG	DA	3342	1/1	0.68	0.94	-	112,112,112,112	0
54	MG	BA	3580	1/1	0.86	0.64	-	94,94,94,94	0
54	MG	BA	3183	1/1	0.76	0.17	-	68,68,68,68	0
54	MG	CL	201	1/1	0.84	0.14	-	99,99,99,99	0
54	MG	BB	227	1/1	0.90	0.43	-	80,80,80,80	0
54	MG	BA	3158	1/1	0.81	0.96	-	99,99,99,99	0
54	MG	DA	3780	1/1	0.78	0.30	-	67,67,67,67	0
54	MG	CA	1669	1/1	0.93	0.17	-	76,76,76,76	0
54	MG	AA	1714	1/1	0.84	0.81	-	81,81,81,81	0
54	MG	DA	3749	1/1	0.81	0.74	-	59,59,59,59	0
54	MG	DA	3561	1/1	0.92	0.29	-	85,85,85,85	0
54	MG	DA	3495	1/1	0.90	0.27	-	111,111,111,111	0
54	MG	BA	3142	1/1	0.94	0.22	-	47,47,47,47	0
54	MG	DA	3346	1/1	0.99	0.15	-	68,68,68,68	0
54	MG	DA	3322	1/1	0.96	0.15	-	65,65,65,65	0
54	MG	DA	2920	1/1	0.96	0.29	-	14,14,14,14	0
54	MG	DB	226	1/1	0.83	0.14	-	73,73,73,73	0
54	MG	DA	3167	1/1	0.90	0.45	-	58,58,58,58	0
54	MG	CA	1785	1/1	0.89	0.26	-	76,76,76,76	0
54	MG	DA	3275	1/1	0.95	0.53	-	92,92,92,92	0
54	MG	CA	1650	1/1	0.47	0.28	-	111,111,111,111	0
54	MG	CA	1617	1/1	0.80	0.19	-	104,104,104,104	0
54	MG	CA	1926	1/1	0.98	0.30	-	76,76,76,76	0
54	MG	AA	1805	1/1	0.64	0.51	-	66,66,66,66	0
54	MG	BA	3447	1/1	0.91	0.80	-	39,39,39,39	0
54	MG	AA	1746	1/1	0.95	0.05	-	86,86,86,86	0
54	MG	DA	3763	1/1	0.97	0.26	-	57,57,57,57	0
54	MG	BB	221	1/1	0.84	0.35	-	89,89,89,89	0
54	MG	CA	1964	1/1	0.95	0.21	-	99,99,99,99	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
54	MG	DA	3042	1/1	0.85	0.29	-	81,81,81,81	0
54	MG	D2	101	1/1	0.67	1.19	-	78,78,78,78	0
54	MG	BA	3440	1/1	0.97	0.41	-	23,23,23,23	0
54	MG	DA	3788	1/1	0.94	0.33	-	92,92,92,92	0
54	MG	AA	1697	1/1	0.86	0.81	-	73,73,73,73	0
54	MG	AA	1749	1/1	0.89	0.24	-	90,90,90,90	0
54	MG	BA	3572	1/1	0.94	0.09	-	87,87,87,87	0
54	MG	DA	3012	1/1	0.95	0.28	-	48,48,48,48	0
54	MG	DA	3304	1/1	0.80	0.28	-	85,85,85,85	0
54	MG	AA	1638	1/1	0.96	0.24	-	89,89,89,89	0
54	MG	BA	3012	1/1	0.84	0.89	-	71,71,71,71	0
54	MG	CA	1887	1/1	0.93	0.15	-	75,75,75,75	0
54	MG	DA	3124	1/1	0.98	0.40	-	51,51,51,51	0
54	MG	DA	3467	1/1	0.99	0.14	-	97,97,97,97	0
54	MG	DA	3818	1/1	0.89	0.16	-	78,78,78,78	0
54	MG	DA	3330	1/1	0.75	0.40	-	100,100,100,100	0
54	MG	CB	303	1/1	0.80	0.10	-	122,122,122,122	0
54	MG	CA	1972	1/1	0.93	0.29	-	67,67,67,67	0
54	MG	BA	3659	1/1	0.86	0.28	-	68,68,68,68	0
54	MG	CA	1750	1/1	0.90	0.27	-	87,87,87,87	0
54	MG	DA	3347	1/1	0.90	0.58	-	67,67,67,67	0
54	MG	AA	1863	1/1	0.94	0.15	-	119,119,119,119	0
54	MG	DA	2952	1/1	0.89	0.47	-	25,25,25,25	0
54	MG	BA	3585	1/1	0.76	0.51	-	69,69,69,69	0
54	MG	DA	3621	1/1	0.73	0.20	-	88,88,88,88	0
54	MG	BA	3323	1/1	0.68	0.32	-	98,98,98,98	0
54	MG	BA	3545	1/1	0.80	0.60	-	59,59,59,59	0
54	MG	CA	1815	1/1	0.92	0.39	-	97,97,97,97	0
54	MG	DK	202	1/1	0.73	0.28	-	69,69,69,69	0
54	MG	DA	3408	1/1	0.88	0.23	-	109,109,109,109	0
54	MG	CA	1839	1/1	0.86	0.35	-	92,92,92,92	0
54	MG	BA	3367	1/1	0.93	0.22	-	74,74,74,74	0
54	MG	BA	3478	1/1	0.85	0.77	-	78,78,78,78	0
54	MG	DA	3505	1/1	0.79	0.40	-	105,105,105,105	0
54	MG	DA	3795	1/1	0.83	0.48	-	80,80,80,80	0
54	MG	BA	2995	1/1	0.76	0.84	-	57,57,57,57	0
54	MG	BA	3251	1/1	0.90	0.54	-	95,95,95,95	0
54	MG	CA	1730	1/1	0.72	0.19	-	112,112,112,112	0
54	MG	BA	3141	1/1	0.90	0.65	-	91,91,91,91	0
54	MG	AA	1744	1/1	0.70	0.17	-	96,96,96,96	0
54	MG	BA	3303	1/1	0.94	0.22	-	157,157,157,157	0
54	MG	DA	2994	1/1	0.95	0.33	-	54,54,54,54	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
54	MG	CA	1777	1/1	0.97	0.34	-	98,98,98,98	0
54	MG	DA	3816	1/1	0.93	0.19	-	52,52,52,52	0
54	MG	AA	1656	1/1	0.83	0.57	-	86,86,86,86	0
54	MG	DA	2992	1/1	0.95	0.39	-	60,60,60,60	0
54	MG	DA	3453	1/1	0.86	0.63	-	87,87,87,87	0
54	MG	DA	3314	1/1	0.85	0.37	-	105,105,105,105	0
54	MG	CA	1800	1/1	0.86	0.18	-	89,89,89,89	0
54	MG	BA	3629	1/1	0.81	0.34	-	86,86,86,86	0
54	MG	DA	3231	1/1	0.77	0.48	-	81,81,81,81	0
54	MG	DA	3113	1/1	0.84	0.16	-	66,66,66,66	0
54	MG	DA	3345	1/1	0.90	0.28	-	88,88,88,88	0
54	MG	AA	1673	1/1	0.92	0.28	-	95,95,95,95	0
54	MG	BO	201	1/1	0.91	0.13	-	70,70,70,70	0
54	MG	BA	3628	1/1	0.94	0.10	-	102,102,102,102	0
54	MG	BA	3287	1/1	0.77	0.33	-	83,83,83,83	0
54	MG	AA	1758	1/1	0.95	0.62	-	69,69,69,69	0
54	MG	BA	3493	1/1	0.79	0.48	-	75,75,75,75	0
54	MG	DA	3150	1/1	0.99	0.12	-	56,56,56,56	0
54	MG	DA	3482	1/1	0.93	0.18	-	69,69,69,69	0
54	MG	DA	3166	1/1	0.86	0.09	-	73,73,73,73	0
54	MG	DA	3240	1/1	0.95	0.16	-	69,69,69,69	0
54	MG	CA	1778	1/1	0.78	0.32	-	94,94,94,94	0
54	MG	DA	3364	1/1	0.94	0.52	-	75,75,75,75	0
54	MG	BJ	203	1/1	0.73	0.23	-	89,89,89,89	0
54	MG	AA	1622	1/1	0.88	0.54	-	62,62,62,62	0
54	MG	DA	3768	1/1	0.73	0.48	-	102,102,102,102	0
54	MG	DA	3738	1/1	0.92	0.20	-	91,91,91,91	0
54	MG	DA	3777	1/1	0.98	0.29	-	90,90,90,90	0
54	MG	BA	2999	1/1	0.96	0.53	-	39,39,39,39	0
54	MG	BK	201	1/1	0.59	0.19	-	77,77,77,77	0
54	MG	BA	3479	1/1	0.82	0.25	-	69,69,69,69	0
54	MG	DA	3465	1/1	0.89	0.27	-	105,105,105,105	0
54	MG	BA	3428	1/1	0.97	0.26	-	90,90,90,90	0
54	MG	CA	1631	1/1	0.91	0.54	-	82,82,82,82	0
54	MG	DJ	202	1/1	0.86	0.37	-	126,126,126,126	0
54	MG	DA	3700	1/1	0.66	1.31	-	96,96,96,96	0
54	MG	BA	3050	1/1	0.88	0.57	-	57,57,57,57	0
54	MG	BA	3061	1/1	0.97	0.13	-	54,54,54,54	0
54	MG	DA	3753	1/1	0.88	0.38	-	79,79,79,79	0
54	MG	AA	1853	1/1	0.82	0.10	-	86,86,86,86	0
54	MG	BA	3473	1/1	0.91	1.49	-	70,70,70,70	0
54	MG	BA	3450	1/1	0.94	0.64	-	46,46,46,46	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
54	MG	BA	3424	1/1	0.98	0.33	-	115,115,115,115	0
54	MG	CV	6202	1/1	0.82	0.12	-	108,108,108,108	0
54	MG	DA	3003	1/1	0.90	0.56	-	51,51,51,51	0
54	MG	BA	3330	1/1	0.84	0.60	-	101,101,101,101	0
54	MG	BB	225	1/1	0.82	1.22	-	96,96,96,96	0
54	MG	BA	3391	1/1	0.91	0.36	-	93,93,93,93	0
54	MG	DA	2965	1/1	0.95	0.60	-	34,34,34,34	0
54	MG	DJ	201	1/1	0.93	0.30	-	61,61,61,61	0
54	MG	CA	1768	1/1	0.87	0.20	-	123,123,123,123	0
54	MG	DA	2989	1/1	0.83	0.36	-	39,39,39,39	0
54	MG	DA	3604	1/1	0.91	0.16	-	103,103,103,103	0
54	MG	AA	1817	1/1	0.92	0.83	-	84,84,84,84	0
54	MG	BA	3437	1/1	0.98	0.53	-	42,42,42,42	0
54	MG	CA	1772	1/1	0.96	0.14	-	112,112,112,112	0
54	MG	DA	3686	1/1	0.90	0.18	-	52,52,52,52	0
54	MG	DA	3129	1/1	0.93	0.45	-	93,93,93,93	0
54	MG	DA	3255	1/1	0.76	0.44	-	70,70,70,70	0
54	MG	DA	3025	1/1	0.81	0.56	-	80,80,80,80	0
54	MG	BA	3337	1/1	0.95	0.18	-	77,77,77,77	0
54	MG	BA	3655	1/1	0.96	0.17	-	92,92,92,92	0
54	MG	DA	2951	1/1	0.89	0.67	-	56,56,56,56	0
54	MG	DA	3558	1/1	0.87	0.27	-	148,148,148,148	0
54	MG	CA	1655	1/1	0.98	0.26	-	100,100,100,100	0
54	MG	DO	201	1/1	0.82	0.53	-	55,55,55,55	0
54	MG	AA	1777	1/1	0.91	0.30	-	49,49,49,49	0
54	MG	DA	3034	1/1	0.77	0.53	-	89,89,89,89	0
54	MG	BA	3489	1/1	0.90	0.18	-	73,73,73,73	0
54	MG	BA	3495	1/1	0.92	0.11	-	74,74,74,74	0
54	MG	DA	3718	1/1	0.94	0.22	-	70,70,70,70	0
54	MG	CA	1653	1/1	0.88	0.16	-	133,133,133,133	0
54	MG	DA	3496	1/1	0.84	0.38	-	95,95,95,95	0
54	MG	BA	3280	1/1	0.79	0.42	-	61,61,61,61	0
54	MG	BA	3219	1/1	0.97	0.15	-	80,80,80,80	0
54	MG	BA	3338	1/1	0.94	0.19	-	78,78,78,78	0
54	MG	BA	3248	1/1	0.87	0.52	-	82,82,82,82	0
54	MG	BA	3624	1/1	0.93	0.12	-	82,82,82,82	0
54	MG	DB	231	1/1	0.86	0.33	-	117,117,117,117	0
54	MG	DA	3220	1/1	0.94	0.17	-	98,98,98,98	0
54	MG	CA	1932	1/1	0.79	0.13	-	84,84,84,84	0
54	MG	BA	3571	1/1	0.96	0.15	-	74,74,74,74	0
54	MG	DA	3540	1/1	0.86	0.21	-	93,93,93,93	0
54	MG	CA	1956	1/1	0.89	0.27	-	85,85,85,85	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
54	MG	BA	3198	1/1	0.67	0.38	-	85,85,85,85	0
54	MG	BA	3188	1/1	0.95	0.19	-	103,103,103,103	0
54	MG	BA	3365	1/1	0.94	0.12	-	131,131,131,131	0
54	MG	DA	3595	1/1	0.91	0.20	-	61,61,61,61	0
54	MG	AN	102	1/1	0.76	0.83	-	91,91,91,91	0
54	MG	BA	3350	1/1	0.86	0.38	-	66,66,66,66	0
54	MG	BA	3062	1/1	0.89	0.53	-	79,79,79,79	0
54	MG	CA	1904	1/1	0.88	0.32	-	97,97,97,97	0
54	MG	AA	1723	1/1	0.90	0.26	-	126,126,126,126	0
54	MG	DA	2933	1/1	0.97	0.32	-	25,25,25,25	0
54	MG	CA	1632	1/1	0.88	0.46	-	59,59,59,59	0
54	MG	AA	1875	1/1	0.88	0.22	-	97,97,97,97	0
54	MG	DA	3734	1/1	0.92	0.38	-	76,76,76,76	0
54	MG	BA	3017	1/1	0.89	0.22	-	55,55,55,55	0
54	MG	DA	3681	1/1	0.96	0.23	-	78,78,78,78	0
54	MG	DA	3259	1/1	0.79	0.15	-	87,87,87,87	0
54	MG	BA	3119	1/1	0.92	0.63	-	73,73,73,73	0
54	MG	BA	2937	1/1	0.90	0.13	-	56,56,56,56	0
54	MG	AA	1733	1/1	0.93	0.20	-	89,89,89,89	0
54	MG	DA	2931	1/1	0.98	0.40	-	21,21,21,21	0
54	MG	DC	303	1/1	0.87	0.88	-	80,80,80,80	0
54	MG	DA	3299	1/1	0.89	0.40	-	104,104,104,104	0
54	MG	CA	1828	1/1	0.91	0.39	-	88,88,88,88	0
54	MG	BA	3452	1/1	0.99	0.25	-	41,41,41,41	0
54	MG	DA	3423	1/1	0.92	0.28	-	79,79,79,79	0
54	MG	DA	3373	1/1	0.87	0.10	-	78,78,78,78	0
54	MG	DA	3102	1/1	0.86	0.26	-	50,50,50,50	0
54	MG	DA	3562	1/1	0.94	0.20	-	60,60,60,60	0
54	MG	DA	3483	1/1	0.98	0.45	-	63,63,63,63	0
54	MG	AA	1867	1/1	0.93	0.18	-	114,114,114,114	0
54	MG	BA	3400	1/1	0.77	0.55	-	83,83,83,83	0
54	MG	DA	3298	1/1	0.89	1.05	-	72,72,72,72	0
54	MG	DA	3140	1/1	0.95	0.73	-	50,50,50,50	0
54	MG	BA	3382	1/1	0.79	0.14	-	85,85,85,85	0
54	MG	DA	3274	1/1	0.87	0.13	-	116,116,116,116	0
54	MG	DA	3574	1/1	0.87	0.26	-	152,152,152,152	0
54	MG	AA	1715	1/1	0.90	0.23	-	58,58,58,58	0
54	MG	DA	3199	1/1	0.97	0.09	-	108,108,108,108	0
54	MG	DB	243	1/1	0.62	0.35	-	101,101,101,101	0
54	MG	CA	1786	1/1	0.92	0.33	-	94,94,94,94	0
54	MG	DA	3703	1/1	0.88	0.37	-	60,60,60,60	0
54	MG	BA	3211	1/1	0.74	0.42	-	100,100,100,100	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
54	MG	CA	1930	1/1	0.90	0.10	-	89,89,89,89	0
54	MG	AA	1708	1/1	0.84	0.37	-	137,137,137,137	0
54	MG	DA	3444	1/1	0.88	0.38	-	72,72,72,72	0
54	MG	DA	3761	1/1	0.96	0.11	-	100,100,100,100	0
54	MG	CA	1677	1/1	0.95	0.45	-	93,93,93,93	0
54	MG	AA	1757	1/1	0.91	0.14	-	98,98,98,98	0
54	MG	DA	3601	1/1	0.95	0.20	-	59,59,59,59	0
54	MG	DA	3401	1/1	0.94	0.31	-	100,100,100,100	0
54	MG	DG	201	1/1	0.79	0.19	-	108,108,108,108	0
54	MG	BA	3096	1/1	0.88	0.33	-	76,76,76,76	0
54	MG	CA	1693	1/1	0.95	0.87	-	95,95,95,95	0
54	MG	CA	1856	1/1	0.83	1.38	-	118,118,118,118	0
54	MG	DA	3740	1/1	0.88	0.10	-	99,99,99,99	0
54	MG	DA	3387	1/1	0.94	0.12	-	124,124,124,124	0
54	MG	BA	3599	1/1	0.44	0.25	-	94,94,94,94	0
54	MG	DA	3842	1/1	0.84	0.37	-	105,105,105,105	0
54	MG	BA	3306	1/1	0.86	0.25	-	78,78,78,78	0
54	MG	CA	1711	1/1	0.84	0.48	-	76,76,76,76	0
54	MG	DA	3217	1/1	0.85	0.31	-	59,59,59,59	0
54	MG	DA	3033	1/1	0.88	0.51	-	79,79,79,79	0
54	MG	AA	1854	1/1	0.60	0.56	-	114,114,114,114	0
54	MG	BA	3086	1/1	0.94	0.11	-	63,63,63,63	0
54	MG	BA	3378	1/1	0.77	0.47	-	80,80,80,80	0
54	MG	BA	3325	1/1	0.99	0.22	-	168,168,168,168	0
54	MG	DA	3789	1/1	0.97	0.15	-	56,56,56,56	0
54	MG	CA	1793	1/1	0.96	0.19	-	125,125,125,125	0
54	MG	DA	3460	1/1	0.96	0.45	-	86,86,86,86	0
54	MG	CA	1883	1/1	0.88	0.38	-	75,75,75,75	0
54	MG	DA	3478	1/1	0.95	0.33	-	48,48,48,48	0
54	MG	AH	201	1/1	0.83	0.28	-	78,78,78,78	0
54	MG	BA	3492	1/1	0.96	0.17	-	70,70,70,70	0
54	MG	AA	1640	1/1	0.94	0.29	-	58,58,58,58	0
54	MG	DA	3790	1/1	0.94	0.48	-	63,63,63,63	0
54	MG	DL	201	1/1	0.98	0.10	-	7,7,7,7	0
54	MG	BA	3137	1/1	0.94	0.29	-	90,90,90,90	0
54	MG	BA	3568	1/1	0.93	0.35	-	69,69,69,69	0
54	MG	BA	3619	1/1	0.98	0.20	-	145,145,145,145	0
54	MG	AA	1850	1/1	0.94	0.70	-	70,70,70,70	0
54	MG	DA	3210	1/1	0.79	0.18	-	92,92,92,92	0
54	MG	BA	3208	1/1	0.83	0.45	-	96,96,96,96	0
54	MG	DE	302	1/1	0.77	0.41	-	83,83,83,83	0
54	MG	CA	1941	1/1	0.81	0.19	-	70,70,70,70	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
54	MG	DA	2916	1/1	0.97	0.47	-	15,15,15,15	0
54	MG	CA	1910	1/1	0.95	0.20	-	109,109,109,109	0
54	MG	AA	1858	1/1	0.89	0.08	-	86,86,86,86	0
54	MG	DA	3779	1/1	0.91	0.23	-	81,81,81,81	0
54	MG	AG	201	1/1	0.82	0.09	-	106,106,106,106	0
54	MG	AA	1732	1/1	0.96	0.13	-	104,104,104,104	0
54	MG	BA	3384	1/1	0.83	0.30	-	93,93,93,93	0
54	MG	DA	3162	1/1	0.90	0.38	-	30,30,30,30	0
54	MG	DA	3292	1/1	0.93	1.15	-	73,73,73,73	0
54	MG	DA	3261	1/1	0.95	0.29	-	57,57,57,57	0
54	MG	DA	3583	1/1	0.83	0.49	-	119,119,119,119	0
54	MG	AA	1833	1/1	0.94	0.43	-	75,75,75,75	0
54	MG	BA	2974	1/1	0.89	0.31	-	81,81,81,81	0
54	MG	DA	3584	1/1	0.88	0.46	-	130,130,130,130	0
54	MG	BU	201	1/1	0.91	0.36	-	70,70,70,70	0
54	MG	AO	103	1/1	0.93	0.10	-	111,111,111,111	0
54	MG	BA	3304	1/1	0.90	0.58	-	100,100,100,100	0
54	MG	CA	1701	1/1	0.91	0.70	-	83,83,83,83	0
54	MG	CA	1717	1/1	0.73	0.67	-	103,103,103,103	0
54	MG	BA	3226	1/1	0.82	0.36	-	59,59,59,59	0
54	MG	DA	3399	1/1	0.95	0.17	-	75,75,75,75	0
54	MG	AA	1731	1/1	0.85	0.32	-	83,83,83,83	0
54	MG	DA	3079	1/1	0.88	0.35	-	51,51,51,51	0
54	MG	DA	3207	1/1	0.89	0.29	-	43,43,43,43	0
54	MG	DA	3077	1/1	0.98	0.09	-	33,33,33,33	0
54	MG	DA	3409	1/1	0.66	0.24	-	110,110,110,110	0
54	MG	DA	3377	1/1	0.78	0.31	-	107,107,107,107	0
54	MG	BA	2910	1/1	0.92	0.39	-	17,17,17,17	0
54	MG	DA	2958	1/1	0.95	0.07	-	31,31,31,31	0
54	MG	CA	1935	1/1	0.88	0.26	-	101,101,101,101	0
54	MG	CA	1939	1/1	0.85	0.07	-	144,144,144,144	0
54	MG	BA	3410	1/1	0.87	0.46	-	112,112,112,112	0
54	MG	DA	3280	1/1	0.84	0.15	-	85,85,85,85	0
54	MG	DB	213	1/1	0.70	0.17	-	113,113,113,113	0
54	MG	DA	3834	1/1	0.98	0.13	-	60,60,60,60	0
54	MG	DA	3692	1/1	0.74	0.68	-	86,86,86,86	0
54	MG	CA	1678	1/1	0.66	0.39	-	81,81,81,81	0
54	MG	CA	1738	1/1	0.84	0.23	-	93,93,93,93	0
54	MG	BA	3637	1/1	0.95	0.18	-	105,105,105,105	0
54	MG	CA	1905	1/1	0.98	0.28	-	130,130,130,130	0
54	MG	CA	1885	1/1	0.83	0.37	-	75,75,75,75	0
54	MG	AA	1679	1/1	0.99	0.34	-	78,78,78,78	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
54	MG	AA	1786	1/1	0.87	0.65	-	119,119,119,119	0
54	MG	DA	3100	1/1	0.88	0.14	-	78,78,78,78	0
54	MG	AA	1619	1/1	0.85	0.27	-	57,57,57,57	0
54	MG	CA	1898	1/1	0.75	0.85	-	87,87,87,87	0
54	MG	DA	3277	1/1	0.75	0.27	-	59,59,59,59	0
54	MG	DA	3557	1/1	0.94	0.12	-	128,128,128,128	0
54	MG	DA	3532	1/1	0.70	0.33	-	106,106,106,106	0
54	MG	CC	306	1/1	0.71	0.40	-	117,117,117,117	0
54	MG	DA	3762	1/1	0.69	0.78	-	103,103,103,103	0
54	MG	DA	3608	1/1	0.87	0.33	-	128,128,128,128	0
54	MG	CA	1648	1/1	0.93	0.43	-	61,61,61,61	0
54	MG	DA	3504	1/1	0.92	0.44	-	97,97,97,97	0
54	MG	DA	3600	1/1	0.63	0.07	-	86,86,86,86	0
54	MG	CA	1805	1/1	0.98	0.09	-	94,94,94,94	0
54	MG	DA	3805	1/1	0.89	0.10	-	70,70,70,70	0
54	MG	DA	3645	1/1	0.97	0.28	-	23,23,23,23	0
54	MG	CA	1608	1/1	0.98	0.35	-	121,121,121,121	0
54	MG	DA	3486	1/1	0.94	0.54	-	124,124,124,124	0
54	MG	BA	3385	1/1	0.29	0.20	-	125,125,125,125	0
54	MG	DA	3778	1/1	0.89	0.52	-	80,80,80,80	0
54	MG	BA	3103	1/1	0.94	0.87	-	74,74,74,74	0
54	MG	DA	3308	1/1	0.91	0.15	-	83,83,83,83	0
54	MG	BA	3117	1/1	0.76	0.47	-	91,91,91,91	0
54	MG	DA	3515	1/1	0.89	0.51	-	73,73,73,73	0
54	MG	CA	1656	1/1	0.97	0.05	-	78,78,78,78	0
54	MG	AA	1840	1/1	0.99	0.05	-	91,91,91,91	0
54	MG	DA	3690	1/1	0.95	0.51	-	70,70,70,70	0
54	MG	D2	103	1/1	0.65	0.51	-	56,56,56,56	0
54	MG	DA	3351	1/1	0.89	0.30	-	89,89,89,89	0
54	MG	DA	3011	1/1	0.98	0.12	-	36,36,36,36	0
54	MG	BA	3203	1/1	0.85	0.14	-	67,67,67,67	0
54	MG	DA	3341	1/1	0.91	0.24	-	75,75,75,75	0
54	MG	DA	3543	1/1	0.91	0.20	-	87,87,87,87	0
54	MG	BA	3084	1/1	0.96	0.37	-	82,82,82,82	0
54	MG	DA	3588	1/1	0.33	0.42	-	96,96,96,96	0
54	MG	CA	1833	1/1	0.89	0.28	-	79,79,79,79	0
54	MG	DA	3030	1/1	0.86	0.28	-	73,73,73,73	0
54	MG	DA	3061	1/1	0.83	0.24	-	55,55,55,55	0
54	MG	DA	3238	1/1	0.88	0.33	-	56,56,56,56	0
54	MG	DA	3454	1/1	0.95	0.36	-	81,81,81,81	0
54	MG	BA	3612	1/1	0.82	0.35	-	85,85,85,85	0
54	MG	BA	3647	1/1	0.86	0.62	-	149,149,149,149	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
54	MG	DA	3725	1/1	0.90	0.45	-	65,65,65,65	0
54	MG	CA	1726	1/1	0.96	0.29	-	49,49,49,49	0
54	MG	BA	3265	1/1	0.86	0.16	-	53,53,53,53	0
54	MG	BA	3253	1/1	0.87	0.18	-	88,88,88,88	0
54	MG	DY	101	1/1	0.85	0.30	-	12,12,12,12	0
54	MG	DB	209	1/1	0.83	0.33	-	103,103,103,103	0
54	MG	BA	3266	1/1	0.90	0.13	-	104,104,104,104	0
54	MG	CA	1744	1/1	0.97	0.24	-	80,80,80,80	0
54	MG	DA	3811	1/1	0.94	0.17	-	60,60,60,60	0
54	MG	BA	3484	1/1	0.93	0.20	-	56,56,56,56	0
54	MG	DY	102	1/1	0.88	0.38	-	69,69,69,69	0
54	MG	DA	3067	1/1	0.74	0.69	-	91,91,91,91	0
54	MG	BA	3368	1/1	0.91	0.07	-	58,58,58,58	0
54	MG	BA	3622	1/1	0.94	0.27	-	90,90,90,90	0
54	MG	CA	1951	1/1	0.39	0.21	-	87,87,87,87	0
54	MG	CA	1666	1/1	0.75	0.33	-	104,104,104,104	0
54	MG	DA	3323	1/1	0.68	1.21	-	78,78,78,78	0
54	MG	DA	3552	1/1	0.93	0.17	-	68,68,68,68	0
54	MG	AA	1752	1/1	0.88	0.24	-	94,94,94,94	0
54	MG	BA	3328	1/1	0.92	0.28	-	80,80,80,80	0
54	MG	DA	3354	1/1	0.85	0.50	-	79,79,79,79	0
54	MG	DA	3477	1/1	0.94	0.16	-	78,78,78,78	0
54	MG	DA	3616	1/1	0.85	0.46	-	103,103,103,103	0
54	MG	BA	3278	1/1	0.93	0.28	-	88,88,88,88	0
54	MG	DA	3356	1/1	0.91	0.17	-	70,70,70,70	0
54	MG	CA	1814	1/1	0.94	0.33	-	47,47,47,47	0
54	MG	BA	3249	1/1	0.91	0.26	-	36,36,36,36	0
54	MG	DA	3572	1/1	0.94	0.14	-	93,93,93,93	0
54	MG	DA	3131	1/1	0.96	0.10	-	63,63,63,63	0
54	MG	CA	1888	1/1	0.61	0.13	-	109,109,109,109	0
54	MG	DA	3317	1/1	0.84	0.29	-	96,96,96,96	0
54	MG	CA	1804	1/1	0.92	2.12	-	100,100,100,100	0
54	MG	BA	3632	1/1	0.93	0.57	-	61,61,61,61	0
54	MG	DA	3652	1/1	0.75	0.58	-	66,66,66,66	0
54	MG	BB	204	1/1	0.96	0.33	-	102,102,102,102	0
54	MG	CA	1963	1/1	0.94	0.17	-	99,99,99,99	0
54	MG	BA	3089	1/1	0.91	0.49	-	69,69,69,69	0
54	MG	DA	3462	1/1	0.79	0.39	-	83,83,83,83	0
54	MG	CA	1649	1/1	0.89	1.00	-	96,96,96,96	0
54	MG	AA	1727	1/1	0.97	0.37	-	104,104,104,104	0
54	MG	BA	3170	1/1	0.96	0.34	-	87,87,87,87	0
54	MG	BA	3156	1/1	0.91	0.42	-	70,70,70,70	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
54	MG	BA	3043	1/1	0.75	0.32	-	50,50,50,50	0
54	MG	CA	1691	1/1	0.94	0.07	-	92,92,92,92	0
54	MG	BA	3133	1/1	0.77	0.41	-	86,86,86,86	0
54	MG	DA	3679	1/1	0.86	0.34	-	50,50,50,50	0
54	MG	BA	3190	1/1	0.86	0.35	-	72,72,72,72	0
54	MG	DA	3190	1/1	0.94	0.43	-	69,69,69,69	0
54	MG	CA	1716	1/1	0.96	0.20	-	89,89,89,89	0
54	MG	BA	3336	1/1	0.89	0.73	-	87,87,87,87	0
54	MG	DA	3378	1/1	0.90	0.36	-	68,68,68,68	0
54	MG	DA	3674	1/1	0.97	0.54	-	47,47,47,47	0
54	MG	DA	3135	1/1	0.88	0.13	-	114,114,114,114	0
54	MG	DJ	203	1/1	0.92	0.84	-	84,84,84,84	0
54	MG	DA	3548	1/1	0.96	0.86	-	110,110,110,110	0
54	MG	DA	3837	1/1	0.74	0.35	-	95,95,95,95	0
54	MG	DA	3847	1/1	0.90	0.28	-	91,91,91,91	0
54	MG	CA	1639	1/1	0.96	0.38	-	59,59,59,59	0
54	MG	DA	3843	1/1	0.92	0.60	-	55,55,55,55	0
54	MG	BA	3153	1/1	0.82	0.24	-	79,79,79,79	0
54	MG	DA	3720	1/1	0.85	0.40	-	79,79,79,79	0
54	MG	AA	1717	1/1	0.32	1.41	-	103,103,103,103	0
54	MG	DA	3594	1/1	0.71	0.20	-	58,58,58,58	0
54	MG	BS	201	1/1	0.94	0.20	-	52,52,52,52	0
54	MG	DA	3164	1/1	0.98	0.22	-	83,83,83,83	0
54	MG	BA	3494	1/1	0.93	0.39	-	66,66,66,66	0
54	MG	BA	3466	1/1	0.83	0.52	-	61,61,61,61	0
54	MG	CA	1708	1/1	0.95	0.13	-	60,60,60,60	0
54	MG	BA	3524	1/1	0.92	0.12	-	107,107,107,107	0
54	MG	BA	3194	1/1	0.85	0.40	-	64,64,64,64	0
54	MG	DA	3559	1/1	0.85	1.17	-	92,92,92,92	0
54	MG	BA	3054	1/1	0.93	0.15	-	65,65,65,65	0
54	MG	CA	1725	1/1	0.85	0.08	-	82,82,82,82	0
54	MG	DE	303	1/1	0.90	0.20	-	81,81,81,81	0
54	MG	DA	3827	1/1	0.83	0.24	-	103,103,103,103	0
54	MG	DA	3198	1/1	0.87	0.31	-	65,65,65,65	0
54	MG	BA	3415	1/1	0.91	0.41	-	128,128,128,128	0
54	MG	CA	1756	1/1	0.88	0.26	-	84,84,84,84	0
54	MG	CA	1795	1/1	0.82	0.20	-	131,131,131,131	0
54	MG	BA	3640	1/1	0.56	0.31	-	97,97,97,97	0
54	MG	BA	3192	1/1	0.91	0.57	-	110,110,110,110	0
54	MG	BA	3162	1/1	0.92	0.08	-	86,86,86,86	0
54	MG	AA	1804	1/1	0.90	1.50	-	80,80,80,80	0
54	MG	AA	1824	1/1	0.75	0.29	-	81,81,81,81	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
54	MG	BA	3041	1/1	0.88	0.36	-	62,62,62,62	0
54	MG	BA	3510	1/1	0.64	0.43	-	86,86,86,86	0
54	MG	BA	3352	1/1	0.93	0.33	-	80,80,80,80	0
54	MG	BA	3342	1/1	0.73	0.37	-	62,62,62,62	0
54	MG	DA	3862	1/1	0.67	0.70	-	119,119,119,119	0
54	MG	AA	1693	1/1	0.80	0.46	-	83,83,83,83	0
54	MG	DA	3581	1/1	0.97	0.12	-	68,68,68,68	0
54	MG	DA	3006	1/1	0.94	0.27	-	60,60,60,60	0
54	MG	DA	3498	1/1	0.98	0.50	-	51,51,51,51	0
54	MG	CA	1858	1/1	0.84	0.87	-	107,107,107,107	0
54	MG	BA	3264	1/1	0.91	0.09	-	87,87,87,87	0
54	MG	AA	1873	1/1	0.86	0.49	-	105,105,105,105	0
54	MG	DU	201	1/1	0.85	0.09	-	84,84,84,84	0
54	MG	DA	3181	1/1	0.86	0.13	-	96,96,96,96	0
54	MG	AA	1604	1/1	0.91	0.46	-	62,62,62,62	0
54	MG	BA	2935	1/1	0.98	0.40	-	26,26,26,26	0
54	MG	D5	102	1/1	0.87	0.24	-	59,59,59,59	0
54	MG	DA	3666	1/1	0.77	0.31	-	81,81,81,81	0
54	MG	BA	3095	1/1	0.54	0.59	-	96,96,96,96	0
54	MG	BA	3186	1/1	0.89	0.20	-	85,85,85,85	0
54	MG	BA	3244	1/1	0.97	0.17	-	79,79,79,79	0
54	MG	BA	3549	1/1	0.65	1.60	-	87,87,87,87	0
54	MG	DA	3568	1/1	0.99	0.10	-	78,78,78,78	0
54	MG	DA	3706	1/1	0.95	0.16	-	75,75,75,75	0
54	MG	CE	201	1/1	0.89	0.13	-	117,117,117,117	0
54	MG	BA	3258	1/1	0.94	0.20	-	78,78,78,78	0
54	MG	DA	3598	1/1	0.83	0.13	-	108,108,108,108	0
54	MG	DA	3838	1/1	0.92	0.48	-	117,117,117,117	0
54	MG	CA	1840	1/1	0.96	0.30	-	76,76,76,76	0
54	MG	DA	3331	1/1	0.94	0.15	-	90,90,90,90	0
54	MG	DA	3573	1/1	0.78	0.43	-	85,85,85,85	0
54	MG	AQ	203	1/1	0.85	0.20	-	101,101,101,101	0
54	MG	AE	205	1/1	0.94	0.43	-	106,106,106,106	0
54	MG	CA	1623	1/1	0.80	0.40	-	93,93,93,93	0
54	MG	DA	3825	1/1	0.90	0.38	-	103,103,103,103	0
54	MG	BA	3071	1/1	0.91	0.21	-	71,71,71,71	0
54	MG	DA	3531	1/1	0.89	0.35	-	93,93,93,93	0
54	MG	DA	3175	1/1	0.87	0.10	-	102,102,102,102	0
54	MG	CA	1728	1/1	0.99	0.15	-	120,120,120,120	0
54	MG	DA	3228	1/1	0.84	0.32	-	54,54,54,54	0
54	MG	BB	217	1/1	0.75	0.34	-	72,72,72,72	0
54	MG	BA	3025	1/1	0.95	0.49	-	76,76,76,76	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
54	MG	DA	3466	1/1	0.94	0.08	-	117,117,117,117	0
54	MG	DA	3828	1/1	0.79	0.24	-	85,85,85,85	0
54	MG	DB	201	1/1	0.93	0.27	-	44,44,44,44	0
54	MG	CA	1813	1/1	0.95	0.17	-	107,107,107,107	0
54	MG	BA	3483	1/1	0.94	0.32	-	71,71,71,71	0
54	MG	DA	3755	1/1	0.86	0.30	-	89,89,89,89	0
54	MG	DA	3204	1/1	0.88	0.50	-	53,53,53,53	0
54	MG	AA	1755	1/1	0.77	0.16	-	98,98,98,98	0
54	MG	CA	1873	1/1	0.92	0.44	-	58,58,58,58	0
54	MG	AA	1695	1/1	0.85	0.12	-	96,96,96,96	0
54	MG	BB	212	1/1	0.87	0.23	-	107,107,107,107	0
54	MG	DA	3807	1/1	0.91	0.26	-	61,61,61,61	0
54	MG	DA	3602	1/1	0.89	0.39	-	91,91,91,91	0
54	MG	AA	1718	1/1	0.93	0.21	-	63,63,63,63	0
54	MG	BA	2965	1/1	0.98	0.23	-	51,51,51,51	0
54	MG	DA	2942	1/1	0.93	0.92	-	54,54,54,54	0
54	MG	CA	1969	1/1	0.87	0.60	-	124,124,124,124	0
54	MG	CM	202	1/1	0.16	0.91	-	164,164,164,164	0
54	MG	AA	1809	1/1	0.92	0.75	-	94,94,94,94	0
54	MG	DA	3407	1/1	0.91	0.17	-	108,108,108,108	0
54	MG	BA	3608	1/1	0.81	0.32	-	101,101,101,101	0
54	MG	BA	2918	1/1	0.96	0.61	-	43,43,43,43	0
54	MG	DA	3189	1/1	0.91	0.22	-	73,73,73,73	0
54	MG	BA	3555	1/1	0.84	0.57	-	66,66,66,66	0
54	MG	BA	3480	1/1	0.89	1.07	-	79,79,79,79	0
54	MG	DA	3564	1/1	0.74	0.13	-	77,77,77,77	0
54	MG	DA	3476	1/1	0.71	0.56	-	87,87,87,87	0
54	MG	DA	2919	1/1	0.98	0.33	-	12,12,12,12	0
54	MG	AA	1730	1/1	0.66	0.37	-	107,107,107,107	0
54	MG	CA	1731	1/1	0.98	0.14	-	76,76,76,76	0
54	MG	BA	3002	1/1	0.89	1.37	-	96,96,96,96	0
54	MG	DA	3151	1/1	0.88	0.26	-	76,76,76,76	0
54	MG	AA	1745	1/1	0.99	0.42	-	119,119,119,119	0
54	MG	BA	3429	1/1	0.89	0.38	-	79,79,79,79	0
54	MG	BA	3000	1/1	0.89	0.46	-	49,49,49,49	0
54	MG	BA	2971	1/1	0.84	0.24	-	76,76,76,76	0
54	MG	DA	3812	1/1	0.94	0.22	-	87,87,87,87	0
54	MG	BA	3551	1/1	0.87	0.17	-	95,95,95,95	0
54	MG	DA	2955	1/1	0.99	0.58	-	45,45,45,45	0
54	MG	DA	3767	1/1	0.72	0.33	-	73,73,73,73	0
54	MG	DA	3339	1/1	0.84	0.22	-	81,81,81,81	0
54	MG	DC	301	1/1	0.94	0.73	-	67,67,67,67	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
54	MG	DA	3660	1/1	0.82	0.34	-	65,65,65,65	0
54	MG	CA	1903	1/1	0.67	0.26	-	84,84,84,84	0
54	MG	DA	3861	1/1	0.89	0.30	-	90,90,90,90	0
54	MG	DA	3628	1/1	0.95	0.32	-	56,56,56,56	0
54	MG	BA	3512	1/1	0.85	0.34	-	77,77,77,77	0
54	MG	BA	3530	1/1	0.93	0.28	-	65,65,65,65	0
54	MG	DA	3355	1/1	0.70	0.42	-	76,76,76,76	0
54	MG	AA	1834	1/1	0.81	1.43	-	77,77,77,77	0
54	MG	BA	2932	1/1	0.98	0.86	-	44,44,44,44	0
54	MG	BA	3333	1/1	0.95	0.48	-	54,54,54,54	0
54	MG	BA	3165	1/1	0.92	0.15	-	73,73,73,73	0
54	MG	BA	3542	1/1	0.93	0.18	-	102,102,102,102	0
54	MG	AA	1737	1/1	0.90	0.25	-	111,111,111,111	0
54	MG	DA	3752	1/1	0.82	0.39	-	74,74,74,74	0
54	MG	DA	2962	1/1	0.87	0.19	-	37,37,37,37	0
54	MG	DA	2979	1/1	0.99	0.20	-	16,16,16,16	0
54	MG	BA	3373	1/1	0.82	0.36	-	61,61,61,61	0
54	MG	BA	3173	1/1	0.86	0.51	-	87,87,87,87	0
54	MG	AA	1826	1/1	0.96	0.38	-	75,75,75,75	0
54	MG	DA	3663	1/1	0.71	0.46	-	75,75,75,75	0
54	MG	DA	3724	1/1	0.91	0.25	-	109,109,109,109	0
54	MG	DA	3848	1/1	0.81	0.96	-	105,105,105,105	0
54	MG	DA	3337	1/1	0.75	0.37	-	88,88,88,88	0
54	MG	DA	3375	1/1	0.91	0.30	-	63,63,63,63	0
54	MG	BA	3486	1/1	0.95	0.20	-	60,60,60,60	0
54	MG	DA	3276	1/1	0.95	0.56	-	86,86,86,86	0
54	MG	DA	3027	1/1	0.97	0.28	-	53,53,53,53	0
54	MG	CA	1787	1/1	0.97	0.76	-	37,37,37,37	0
54	MG	BA	3560	1/1	0.95	0.62	-	74,74,74,74	0
54	MG	DA	3194	1/1	0.93	0.44	-	74,74,74,74	0
54	MG	DA	3271	1/1	0.68	0.35	-	99,99,99,99	0
54	MG	BA	3639	1/1	0.97	0.23	-	88,88,88,88	0
54	MG	AA	1869	1/1	0.99	0.08	-	74,74,74,74	0
54	MG	BA	3507	1/1	0.80	0.32	-	68,68,68,68	0
54	MG	CA	1746	1/1	0.88	0.24	-	141,141,141,141	0
54	MG	BA	3544	1/1	0.87	0.39	-	81,81,81,81	0
54	MG	BA	3024	1/1	0.95	0.37	-	52,52,52,52	0
54	MG	DA	3719	1/1	0.94	0.33	-	84,84,84,84	0
54	MG	BP	201	1/1	0.92	0.18	-	122,122,122,122	0
54	MG	DA	3246	1/1	0.92	0.12	-	57,57,57,57	0
54	MG	CA	1824	1/1	0.87	0.21	-	123,123,123,123	0
54	MG	BA	3134	1/1	0.75	0.41	-	90,90,90,90	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
54	MG	AA	1661	1/1	0.87	0.16	-	51,51,51,51	0
54	MG	DA	2996	1/1	0.96	0.45	-	32,32,32,32	0
54	MG	DA	3835	1/1	0.94	0.35	-	111,111,111,111	0
54	MG	CA	1739	1/1	0.92	0.15	-	102,102,102,102	0
54	MG	CC	308	1/1	0.79	0.25	-	110,110,110,110	0
54	MG	BA	3206	1/1	0.72	0.23	-	96,96,96,96	0
54	MG	AA	1618	1/1	0.95	0.62	-	71,71,71,71	0
54	MG	BA	3267	1/1	0.90	0.49	-	65,65,65,65	0
54	MG	CG	202	1/1	0.78	0.62	-	131,131,131,131	0
54	MG	DA	3657	1/1	0.83	0.76	-	57,57,57,57	0
54	MG	DA	3334	1/1	0.95	0.30	-	53,53,53,53	0
54	MG	AA	1795	1/1	0.90	0.34	-	111,111,111,111	0
54	MG	DA	3635	1/1	0.94	0.12	-	73,73,73,73	0
54	MG	CA	1637	1/1	0.83	0.09	-	98,98,98,98	0
54	MG	BA	3374	1/1	0.91	0.36	-	108,108,108,108	0
54	MG	DA	2940	1/1	0.96	0.50	-	31,31,31,31	0
54	MG	DA	3695	1/1	0.89	0.28	-	76,76,76,76	0
54	MG	AA	1672	1/1	0.81	0.67	-	78,78,78,78	0
54	MG	BA	3376	1/1	0.85	0.34	-	123,123,123,123	0
54	MG	DA	3002	1/1	0.90	0.16	-	52,52,52,52	0
54	MG	CA	1922	1/1	0.93	0.08	-	127,127,127,127	0
54	MG	DA	3022	1/1	0.88	0.24	-	35,35,35,35	0
54	MG	BA	3152	1/1	0.84	0.12	-	76,76,76,76	0
54	MG	BA	3358	1/1	0.85	0.41	-	81,81,81,81	0
54	MG	AE	204	1/1	0.91	0.23	-	74,74,74,74	0
54	MG	BA	3090	1/1	0.80	0.44	-	84,84,84,84	0
54	MG	BA	3228	1/1	0.97	0.13	-	42,42,42,42	0
54	MG	AA	1764	1/1	0.84	0.62	-	99,99,99,99	0
54	MG	DA	3202	1/1	0.94	0.20	-	76,76,76,76	0
54	MG	AA	1871	1/1	0.65	0.13	-	110,110,110,110	0
54	MG	CE	204	1/1	0.74	0.34	-	109,109,109,109	0
54	MG	DD	306	1/1	0.72	0.18	-	107,107,107,107	0
54	MG	CA	1896	1/1	0.73	0.34	-	70,70,70,70	0
54	MG	CA	1946	1/1	0.94	0.14	-	137,137,137,137	0
54	MG	DA	3417	1/1	0.79	0.65	-	72,72,72,72	0
54	MG	DA	3344	1/1	0.97	0.16	-	75,75,75,75	0
54	MG	CA	1919	1/1	0.68	0.22	-	145,145,145,145	0
54	MG	DA	3363	1/1	0.94	0.17	-	74,74,74,74	0
54	MG	AA	1847	1/1	0.78	0.33	-	90,90,90,90	0
54	MG	AA	1713	1/1	0.98	0.38	-	98,98,98,98	0
54	MG	CV	6205	1/1	0.21	0.11	-	159,159,159,159	0
54	MG	CA	1754	1/1	0.90	0.86	-	121,121,121,121	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
54	MG	BA	3476	1/1	0.85	0.33	-	53,53,53,53	0
54	MG	CA	1660	1/1	0.88	0.14	-	158,158,158,158	0
54	MG	AA	1798	1/1	0.96	0.45	-	46,46,46,46	0
54	MG	DA	3107	1/1	0.95	0.14	-	50,50,50,50	0
54	MG	AO	102	1/1	0.72	0.34	-	114,114,114,114	0
54	MG	AA	1774	1/1	0.92	0.10	-	70,70,70,70	0
54	MG	DA	3624	1/1	0.88	0.27	-	92,92,92,92	0
54	MG	CA	1948	1/1	0.89	0.07	-	99,99,99,99	0
54	MG	AA	1837	1/1	0.94	0.39	-	115,115,115,115	0
54	MG	DA	3501	1/1	0.82	0.25	-	96,96,96,96	0
54	MG	CA	1629	1/1	0.86	0.34	-	93,93,93,93	0
54	MG	AA	1657	1/1	0.93	0.54	-	79,79,79,79	0
54	MG	DA	3359	1/1	0.71	0.31	-	96,96,96,96	0
54	MG	AA	1836	1/1	0.96	0.08	-	91,91,91,91	0
54	MG	AA	1736	1/1	0.94	0.14	-	112,112,112,112	0
54	MG	CA	1933	1/1	0.73	0.24	-	137,137,137,137	0
54	MG	CA	1640	1/1	0.73	1.18	-	133,133,133,133	0
54	MG	BA	3079	1/1	0.97	0.10	-	73,73,73,73	0
54	MG	DA	3435	1/1	0.92	0.33	-	97,97,97,97	0
54	MG	BA	3254	1/1	0.96	0.34	-	68,68,68,68	0
54	MG	BA	3274	1/1	0.70	0.73	-	98,98,98,98	0
54	MG	BA	3116	1/1	0.92	0.17	-	70,70,70,70	0
54	MG	BA	3118	1/1	0.95	0.14	-	95,95,95,95	0
54	MG	DA	3522	1/1	0.81	0.52	-	147,147,147,147	0
54	MG	CG	201	1/1	0.73	0.48	-	104,104,104,104	0
54	MG	DA	3182	1/1	0.83	0.29	-	63,63,63,63	0
54	MG	CA	1806	1/1	0.94	0.18	-	110,110,110,110	0
54	MG	AA	1690	1/1	0.76	0.30	-	88,88,88,88	0
54	MG	CA	1884	1/1	0.77	0.23	-	99,99,99,99	0
54	MG	CA	1705	1/1	0.90	0.37	-	122,122,122,122	0
54	MG	DA	2935	1/1	0.97	0.50	-	23,23,23,23	0
54	MG	BA	2961	1/1	0.98	0.62	-	51,51,51,51	0
54	MG	BA	3508	1/1	0.87	0.15	-	90,90,90,90	0
54	MG	DB	236	1/1	0.72	0.18	-	84,84,84,84	0
54	MG	AA	1653	1/1	0.74	0.07	-	102,102,102,102	0
54	MG	BA	3112	1/1	0.84	0.30	-	65,65,65,65	0
54	MG	CA	1758	1/1	0.89	0.22	-	77,77,77,77	0
54	MG	DA	3497	1/1	0.53	0.73	-	62,62,62,62	0
54	MG	BA	3405	1/1	0.94	0.16	-	68,68,68,68	0
54	MG	DA	3765	1/1	0.94	0.18	-	64,64,64,64	0
54	MG	DA	3775	1/1	0.87	0.21	-	98,98,98,98	0
54	MG	D5	101	1/1	0.92	0.19	-	46,46,46,46	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
54	MG	AK	201	1/1	0.87	0.33	-	101,101,101,101	0
54	MG	BA	3124	1/1	0.77	0.14	-	87,87,87,87	0
54	MG	BA	3579	1/1	0.94	0.73	-	107,107,107,107	0
54	MG	AA	1849	1/1	0.88	0.31	-	79,79,79,79	0
54	MG	DA	3070	1/1	0.83	0.05	-	85,85,85,85	0
54	MG	BA	3353	1/1	0.84	0.47	-	85,85,85,85	0
54	MG	AA	1815	1/1	0.72	1.70	-	78,78,78,78	0
54	MG	DA	3715	1/1	0.90	0.30	-	57,57,57,57	0
54	MG	DA	3288	1/1	0.90	0.15	-	104,104,104,104	0
54	MG	BA	3343	1/1	0.91	0.25	-	87,87,87,87	0
54	MG	BB	226	1/1	0.67	0.71	-	98,98,98,98	0
54	MG	DA	3096	1/1	0.88	0.51	-	59,59,59,59	0
54	MG	CA	1801	1/1	0.95	0.25	-	82,82,82,82	0
54	MG	DA	2968	1/1	0.96	0.63	-	36,36,36,36	0
54	MG	BB	207	1/1	0.97	0.16	-	111,111,111,111	0
54	MG	DA	3735	1/1	0.82	0.17	-	97,97,97,97	0
54	MG	BA	3037	1/1	0.92	0.24	-	62,62,62,62	0
54	MG	AB	303	1/1	0.61	0.26	-	108,108,108,108	0
54	MG	DB	228	1/1	0.84	0.28	-	88,88,88,88	0
54	MG	DB	206	1/1	0.96	0.27	-	63,63,63,63	0
54	MG	BA	3225	1/1	0.96	0.32	-	54,54,54,54	0
54	MG	DA	3782	1/1	0.94	0.42	-	67,67,67,67	0
54	MG	DA	3623	1/1	0.97	0.41	-	104,104,104,104	0
54	MG	B4	101	1/1	0.89	0.28	-	91,91,91,91	0
54	MG	DA	3721	1/1	0.70	0.80	-	88,88,88,88	0
54	MG	AC	302	1/1	0.90	0.33	-	71,71,71,71	0
54	MG	DA	3503	1/1	0.87	0.28	-	91,91,91,91	0
54	MG	BA	3491	1/1	0.93	0.47	-	79,79,79,79	0
54	MG	AA	1868	1/1	0.80	0.28	-	101,101,101,101	0
54	MG	DA	3580	1/1	0.88	0.32	-	162,162,162,162	0
54	MG	BA	3652	1/1	0.83	0.35	-	109,109,109,109	0
54	MG	CC	305	1/1	0.96	0.08	-	93,93,93,93	0
54	MG	BJ	201	1/1	0.79	0.17	-	75,75,75,75	0
54	MG	BH	201	1/1	0.91	0.24	-	107,107,107,107	0
54	MG	AA	1637	1/1	0.93	0.16	-	62,62,62,62	0
54	MG	DA	3004	1/1	0.95	0.21	-	38,38,38,38	0
54	MG	BA	3268	1/1	0.93	0.64	-	84,84,84,84	0
54	MG	DA	3631	1/1	0.99	0.04	-	73,73,73,73	0
54	MG	AA	1812	1/1	0.94	0.93	-	81,81,81,81	0
54	MG	CA	1635	1/1	0.62	0.36	-	99,99,99,99	0
54	MG	CA	1908	1/1	0.73	0.43	-	104,104,104,104	0
54	MG	DA	3221	1/1	0.94	0.23	-	73,73,73,73	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
54	MG	DA	3668	1/1	0.79	0.56	-	133,133,133,133	0
54	MG	AA	1793	1/1	0.88	0.25	-	68,68,68,68	0
54	MG	DA	3116	1/1	0.90	0.47	-	52,52,52,52	0
54	MG	BA	3161	1/1	0.94	0.34	-	64,64,64,64	0
54	MG	CA	1622	1/1	0.61	0.82	-	77,77,77,77	0
54	MG	DA	3815	1/1	0.98	0.53	-	74,74,74,74	0
54	MG	DA	3741	1/1	0.84	0.18	-	84,84,84,84	0
54	MG	BA	2930	1/1	0.93	0.40	-	47,47,47,47	0
54	MG	DB	241	1/1	0.90	0.13	-	90,90,90,90	0
54	MG	DA	3043	1/1	0.96	0.14	-	59,59,59,59	0
54	MG	DA	3570	1/1	0.65	0.59	-	90,90,90,90	0
54	MG	DA	3225	1/1	0.72	0.11	-	139,139,139,139	0
54	MG	BA	3674	1/1	0.95	0.37	-	55,55,55,55	0
54	MG	BA	3664	1/1	0.94	0.10	-	69,69,69,69	0
54	MG	DA	3293	1/1	0.94	0.23	-	66,66,66,66	0
54	MG	AV	6206	1/1	0.84	0.20	-	87,87,87,87	0
54	MG	CA	1853	1/1	0.92	0.18	-	113,113,113,113	0
54	MG	AA	1776	1/1	0.92	0.37	-	108,108,108,108	0
54	MG	BA	3396	1/1	0.95	0.15	-	108,108,108,108	0
54	MG	DA	3171	1/1	0.89	0.14	-	93,93,93,93	0
54	MG	BA	3066	1/1	0.83	1.05	-	91,91,91,91	0
54	MG	DB	242	1/1	0.89	0.09	-	113,113,113,113	0
54	MG	AQ	202	1/1	0.91	0.13	-	85,85,85,85	0
54	MG	DA	3764	1/1	0.90	0.35	-	80,80,80,80	0
54	MG	DA	3797	1/1	0.79	0.26	-	101,101,101,101	0
54	MG	DA	3593	1/1	0.97	0.43	-	125,125,125,125	0
54	MG	BA	3676	1/1	0.95	0.11	-	112,112,112,112	0
54	MG	DA	3864	1/1	0.95	0.37	-	53,53,53,53	0
54	MG	BA	3269	1/1	0.84	0.71	-	104,104,104,104	0
54	MG	DA	3168	1/1	0.93	0.63	-	67,67,67,67	0
54	MG	AA	1743	1/1	0.86	0.53	-	140,140,140,140	0
54	MG	AA	1659	1/1	0.87	0.52	-	74,74,74,74	0
54	MG	CA	1723	1/1	0.84	0.74	-	85,85,85,85	0
54	MG	DA	3132	1/1	0.86	0.57	-	62,62,62,62	0
54	MG	CA	1766	1/1	0.77	0.12	-	77,77,77,77	0
54	MG	BA	3431	1/1	0.97	0.33	-	26,26,26,26	0
54	MG	BA	3235	1/1	0.89	0.26	-	54,54,54,54	0
54	MG	DA	3294	1/1	0.97	0.17	-	36,36,36,36	0
54	MG	BA	3078	1/1	0.87	0.27	-	104,104,104,104	0
54	MG	DA	3413	1/1	0.96	0.23	-	60,60,60,60	0
54	MG	BA	3525	1/1	0.77	0.36	-	67,67,67,67	0
54	MG	DA	3830	1/1	0.89	0.18	-	125,125,125,125	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
54	MG	BA	3215	1/1	0.85	0.65	-	64,64,64,64	0
54	MG	DB	234	1/1	0.57	0.16	-	105,105,105,105	0
54	MG	BA	3261	1/1	0.75	1.47	-	92,92,92,92	0
54	MG	CA	1902	1/1	0.86	0.18	-	77,77,77,77	0
54	MG	BA	3139	1/1	0.94	0.40	-	100,100,100,100	0
54	MG	DA	3673	1/1	0.81	0.20	-	85,85,85,85	0
54	MG	DA	3066	1/1	0.95	0.24	-	114,114,114,114	0
54	MG	BA	3189	1/1	0.82	0.29	-	76,76,76,76	0
54	MG	DA	3400	1/1	0.93	0.11	-	127,127,127,127	0
54	MG	BA	3427	1/1	0.95	0.10	-	74,74,74,74	0
54	MG	CA	1931	1/1	0.94	0.09	-	99,99,99,99	0
54	MG	DA	3516	1/1	0.88	0.33	-	73,73,73,73	0
54	MG	BA	3500	1/1	0.95	0.29	-	72,72,72,72	0
54	MG	DA	3747	1/1	0.92	0.21	-	78,78,78,78	0
54	MG	DA	3615	1/1	0.96	0.20	-	99,99,99,99	0
54	MG	DA	3799	1/1	0.87	0.36	-	96,96,96,96	0
54	MG	BA	3393	1/1	0.63	0.48	-	136,136,136,136	0
54	MG	CO	103	1/1	0.85	0.99	-	203,203,203,203	0
54	MG	CA	1848	1/1	0.96	0.29	-	96,96,96,96	0
54	MG	DA	3256	1/1	0.94	0.44	-	65,65,65,65	0
54	MG	DA	3691	1/1	0.70	0.62	-	96,96,96,96	0
54	MG	CO	102	1/1	0.50	0.49	-	103,103,103,103	0
54	MG	CA	1880	1/1	0.85	1.03	-	83,83,83,83	0
54	MG	DA	3009	1/1	0.93	0.30	-	54,54,54,54	0
54	MG	BA	3083	1/1	0.89	0.08	-	70,70,70,70	0
54	MG	CA	1970	1/1	0.98	0.08	-	98,98,98,98	0
54	MG	CC	302	1/1	0.98	0.20	-	79,79,79,79	0
54	MG	CA	1662	1/1	0.96	0.13	-	84,84,84,84	0
54	MG	DM	1400	1/1	0.82	0.45	-	106,106,106,106	0
54	MG	CA	1909	1/1	0.72	0.23	-	135,135,135,135	0
54	MG	DA	3485	1/1	0.95	0.31	-	119,119,119,119	0
54	MG	BA	3290	1/1	0.89	0.25	-	79,79,79,79	0
54	MG	BA	3085	1/1	0.92	0.34	-	53,53,53,53	0
54	MG	CA	1694	1/1	0.73	0.12	-	79,79,79,79	0
54	MG	DA	3249	1/1	0.93	0.11	-	111,111,111,111	0
54	MG	BA	3490	1/1	0.96	0.09	-	67,67,67,67	0
54	MG	AA	1709	1/1	0.90	0.58	-	87,87,87,87	0
54	MG	DA	3157	1/1	0.98	0.18	-	33,33,33,33	0
54	MG	CA	1612	1/1	0.75	0.34	-	76,76,76,76	0
54	MG	DA	3019	1/1	0.95	0.28	-	48,48,48,48	0
54	MG	BB	209	1/1	0.93	0.17	-	61,61,61,61	0
54	MG	DA	3114	1/1	0.83	0.38	-	66,66,66,66	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
54	MG	CA	1667	1/1	0.94	0.29	-	77,77,77,77	0
54	MG	DA	3232	1/1	0.97	0.13	-	67,67,67,67	0
54	MG	CA	1713	1/1	0.93	0.12	-	92,92,92,92	0
54	MG	DA	3737	1/1	0.93	0.71	-	136,136,136,136	0
54	MG	D4	101	1/1	0.90	0.22	-	55,55,55,55	0
54	MG	DB	233	1/1	0.94	0.42	-	67,67,67,67	0
54	MG	AA	1740	1/1	0.86	0.13	-	76,76,76,76	0
54	MG	CA	1780	1/1	0.86	0.15	-	93,93,93,93	0
54	MG	DA	3319	1/1	0.91	0.30	-	38,38,38,38	0
54	MG	AA	1636	1/1	0.82	0.11	-	96,96,96,96	0
54	MG	BA	3076	1/1	0.89	0.51	-	68,68,68,68	0
54	MG	B3	101	1/1	0.80	0.16	-	96,96,96,96	0
54	MG	BA	3439	1/1	0.96	0.39	-	17,17,17,17	0
54	MG	BA	3419	1/1	0.90	0.30	-	67,67,67,67	0
54	MG	AA	1788	1/1	0.88	0.13	-	148,148,148,148	0
54	MG	DA	3528	1/1	0.98	0.08	-	87,87,87,87	0
54	MG	BA	3667	1/1	0.97	0.16	-	85,85,85,85	0
54	MG	BA	3097	1/1	0.95	0.27	-	57,57,57,57	0
54	MG	DA	3792	1/1	0.93	0.35	-	40,40,40,40	0
54	MG	CA	1714	1/1	0.80	0.36	-	123,123,123,123	0
54	MG	DA	3520	1/1	0.85	0.20	-	133,133,133,133	0
54	MG	DA	3470	1/1	0.95	0.21	-	127,127,127,127	0
54	MG	BA	3067	1/1	0.74	0.49	-	103,103,103,103	0
54	MG	CL	203	1/1	0.58	0.23	-	123,123,123,123	0
54	MG	BA	3120	1/1	0.73	0.11	-	85,85,85,85	0
54	MG	BA	2905	1/1	0.98	0.63	-	19,19,19,19	0
54	MG	BA	3245	1/1	0.95	0.32	-	95,95,95,95	0
54	MG	DA	3101	1/1	0.81	0.46	-	75,75,75,75	0
54	MG	CA	1924	1/1	0.89	0.19	-	116,116,116,116	0
54	MG	AA	1726	1/1	0.87	0.12	-	103,103,103,103	0
54	MG	DA	2961	1/1	0.92	0.42	-	58,58,58,58	0
54	MG	DA	3388	1/1	0.94	0.24	-	51,51,51,51	0
54	MG	BA	3055	1/1	0.75	0.62	-	78,78,78,78	0
54	MG	AA	1790	1/1	0.66	0.12	-	109,109,109,109	0
54	MG	AA	1704	1/1	0.94	0.16	-	77,77,77,77	0
54	MG	BB	205	1/1	0.89	0.14	-	108,108,108,108	0
54	MG	BA	3222	1/1	0.81	0.24	-	86,86,86,86	0
54	MG	AA	1876	1/1	0.78	0.12	-	80,80,80,80	0
54	MG	DA	3801	1/1	0.85	0.30	-	83,83,83,83	0
54	MG	CA	1692	1/1	0.91	1.05	-	115,115,115,115	0
54	MG	DA	3449	1/1	0.95	0.15	-	79,79,79,79	0
54	MG	AA	1652	1/1	0.83	0.33	-	95,95,95,95	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
54	MG	CA	1765	1/1	0.89	0.19	-	95,95,95,95	0
54	MG	CA	1949	1/1	0.64	0.21	-	98,98,98,98	0
54	MG	DA	3853	1/1	0.81	0.28	-	114,114,114,114	0
54	MG	BA	3291	1/1	0.85	0.39	-	102,102,102,102	0
54	MG	BA	3009	1/1	0.90	0.43	-	73,73,73,73	0
54	MG	CA	1918	1/1	0.81	0.68	-	84,84,84,84	0
54	MG	BA	3499	1/1	0.93	0.29	-	66,66,66,66	0
54	MG	BA	3204	1/1	0.76	0.55	-	68,68,68,68	0
54	MG	CA	1962	1/1	0.46	0.19	-	96,96,96,96	0
54	MG	DA	3502	1/1	0.82	0.34	-	93,93,93,93	0
54	MG	BV	201	1/1	0.92	0.27	-	98,98,98,98	0
54	MG	AV	6203	1/1	0.96	0.15	-	62,62,62,62	0
54	MG	BA	2947	1/1	0.93	0.73	-	37,37,37,37	0
54	MG	DA	3856	1/1	0.98	0.49	-	86,86,86,86	0
54	MG	AA	1861	1/1	0.82	0.22	-	81,81,81,81	0
54	MG	BA	2993	1/1	0.88	0.48	-	49,49,49,49	0
54	MG	AA	1772	1/1	0.91	0.31	-	111,111,111,111	0
54	MG	BA	3167	1/1	0.82	0.63	-	87,87,87,87	0
54	MG	DA	3683	1/1	0.87	0.34	-	58,58,58,58	0
54	MG	BA	3606	1/1	0.73	0.68	-	118,118,118,118	0
54	MG	DA	3263	1/1	0.91	0.17	-	55,55,55,55	0
54	MG	B2	102	1/1	0.86	0.72	-	65,65,65,65	0
54	MG	AA	1682	1/1	0.93	0.16	-	91,91,91,91	0
54	MG	BA	3616	1/1	0.84	0.23	-	89,89,89,89	0
54	MG	AA	1691	1/1	0.97	0.11	-	69,69,69,69	0
54	MG	BA	3366	1/1	0.89	0.32	-	80,80,80,80	0
54	MG	BA	3270	1/1	0.89	0.14	-	93,93,93,93	0
54	MG	BA	3390	1/1	0.93	0.09	-	103,103,103,103	0
54	MG	DA	3567	1/1	0.86	1.01	-	87,87,87,87	0
54	MG	DA	3327	1/1	0.85	0.27	-	79,79,79,79	0
54	MG	BA	3140	1/1	0.83	0.34	-	62,62,62,62	0
54	MG	BB	222	1/1	0.96	0.23	-	112,112,112,112	0
54	MG	DA	3802	1/1	0.95	0.14	-	84,84,84,84	0
54	MG	DA	3286	1/1	0.90	0.23	-	76,76,76,76	0
54	MG	BN	202	1/1	0.91	0.66	-	78,78,78,78	0
54	MG	CV	6207	1/1	0.92	0.20	-	94,94,94,94	0
54	MG	DA	3021	1/1	0.70	1.25	-	68,68,68,68	0
54	MG	CA	1955	1/1	0.94	0.28	-	98,98,98,98	0
54	MG	BA	3596	1/1	0.81	0.23	-	71,71,71,71	0
54	MG	AA	1605	1/1	0.92	0.17	-	65,65,65,65	0
54	MG	DA	3817	1/1	0.86	0.40	-	76,76,76,76	0
54	MG	BA	2970	1/1	0.81	0.42	-	55,55,55,55	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
54	MG	CA	1610	1/1	0.94	0.44	-	90,90,90,90	0
54	MG	CE	205	1/1	0.95	0.49	-	104,104,104,104	0
54	MG	BA	3230	1/1	0.93	0.23	-	111,111,111,111	0
54	MG	DA	2949	1/1	0.94	0.37	-	52,52,52,52	0
54	MG	DB	220	1/1	0.71	0.24	-	75,75,75,75	0
54	MG	BA	3059	1/1	0.96	0.64	-	48,48,48,48	0
54	MG	AA	1773	1/1	0.91	0.08	-	97,97,97,97	0
54	MG	AA	1738	1/1	0.74	0.44	-	104,104,104,104	0
54	MG	DB	238	1/1	0.91	0.16	-	74,74,74,74	0
54	MG	DA	3008	1/1	0.95	0.15	-	50,50,50,50	0
54	MG	CA	1769	1/1	0.97	0.13	-	67,67,67,67	0
54	MG	DA	3402	1/1	0.75	0.26	-	95,95,95,95	0
54	MG	DA	3678	1/1	0.93	0.22	-	58,58,58,58	0
54	MG	BA	3297	1/1	0.77	0.68	-	98,98,98,98	0
54	MG	BA	2976	1/1	0.93	0.30	-	65,65,65,65	0
54	MG	DA	3806	1/1	0.97	0.17	-	92,92,92,92	0
54	MG	BA	3607	1/1	0.84	0.76	-	62,62,62,62	0
54	MG	BA	3558	1/1	0.74	0.36	-	91,91,91,91	0
54	MG	AA	1707	1/1	0.92	0.82	-	86,86,86,86	0
54	MG	CA	1682	1/1	0.93	0.54	-	88,88,88,88	0
54	MG	BA	3441	1/1	0.97	0.37	-	35,35,35,35	0
54	MG	CA	1674	1/1	0.91	0.58	-	70,70,70,70	0
54	MG	DA	3524	1/1	0.95	0.12	-	87,87,87,87	0
54	MG	DA	3480	1/1	0.94	0.15	-	76,76,76,76	0
54	MG	DA	3776	1/1	0.97	0.24	-	172,172,172,172	0
54	MG	DA	3826	1/1	0.70	0.17	-	100,100,100,100	0
54	MG	BA	3645	1/1	0.91	0.22	-	69,69,69,69	0
54	MG	CA	1603	1/1	0.92	0.54	-	48,48,48,48	0
54	MG	D4	104	1/1	0.79	0.28	-	76,76,76,76	0
54	MG	CA	1890	1/1	0.86	0.58	-	78,78,78,78	0
54	MG	DA	3254	1/1	0.72	0.67	-	71,71,71,71	0
54	MG	BA	3532	1/1	0.59	0.24	-	107,107,107,107	0
54	MG	DB	230	1/1	0.50	0.68	-	114,114,114,114	0
54	MG	BA	3380	1/1	0.90	0.26	-	55,55,55,55	0
54	MG	BA	3052	1/1	0.49	0.85	-	73,73,73,73	0
54	MG	DA	3234	1/1	0.69	0.36	-	64,64,64,64	0
54	MG	DB	229	1/1	0.69	0.57	-	134,134,134,134	0
54	MG	DA	3858	1/1	0.83	0.32	-	58,58,58,58	0
54	MG	AA	1785	1/1	0.65	0.98	-	117,117,117,117	0
54	MG	BA	3576	1/1	0.77	0.67	-	83,83,83,83	0
54	MG	BA	3513	1/1	0.91	0.16	-	74,74,74,74	0
54	MG	BA	3177	1/1	0.91	0.08	-	119,119,119,119	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
54	MG	BA	3654	1/1	0.86	0.33	-	122,122,122,122	0
54	MG	DA	3713	1/1	0.78	0.56	-	96,96,96,96	0
54	MG	BA	3105	1/1	0.70	0.18	-	72,72,72,72	0
54	MG	BA	3550	1/1	0.92	0.18	-	79,79,79,79	0
54	MG	DA	3272	1/1	0.92	0.56	-	87,87,87,87	0
54	MG	BA	3613	1/1	0.70	0.21	-	103,103,103,103	0
54	MG	DG	203	1/1	0.65	0.20	-	98,98,98,98	0
54	MG	BA	3277	1/1	0.82	0.34	-	110,110,110,110	0
54	MG	BA	3681	1/1	0.80	0.26	-	95,95,95,95	0
54	MG	CL	204	1/1	0.81	0.25	-	86,86,86,86	0
54	MG	CA	1850	1/1	0.91	0.25	-	129,129,129,129	0
54	MG	CA	1763	1/1	0.58	0.09	-	90,90,90,90	0
54	MG	BA	3564	1/1	0.86	0.76	-	146,146,146,146	0
54	MG	DA	3860	1/1	0.71	0.28	-	78,78,78,78	0
54	MG	CA	1942	1/1	0.97	0.20	-	104,104,104,104	0
54	MG	BA	3446	1/1	0.94	0.42	-	53,53,53,53	0
54	MG	BA	3329	1/1	0.92	0.29	-	113,113,113,113	0
54	MG	CA	1673	1/1	0.98	0.49	-	61,61,61,61	0
54	MG	AA	1616	1/1	0.97	0.65	-	54,54,54,54	0
54	MG	AA	1694	1/1	0.95	0.10	-	71,71,71,71	0
54	MG	CA	1960	1/1	0.82	0.15	-	91,91,91,91	0
54	MG	BV	202	1/1	0.73	0.09	-	75,75,75,75	0
54	MG	BA	3260	1/1	0.94	0.35	-	64,64,64,64	0
54	MG	CA	1877	1/1	0.82	0.26	-	126,126,126,126	0
54	MG	AA	1879	1/1	0.88	0.62	-	119,119,119,119	0
54	MG	AA	1780	1/1	0.90	0.06	-	101,101,101,101	0
54	MG	BA	3180	1/1	0.86	0.34	-	88,88,88,88	0
54	MG	BA	2992	1/1	0.93	0.13	-	69,69,69,69	0
54	MG	DA	3336	1/1	0.89	1.01	-	70,70,70,70	0
54	MG	DA	3247	1/1	0.93	0.14	-	92,92,92,92	0
54	MG	CA	1917	1/1	0.77	0.30	-	105,105,105,105	0
54	MG	DA	3739	1/1	0.98	0.18	-	133,133,133,133	0
54	MG	BA	3592	1/1	0.63	0.53	-	79,79,79,79	0
54	MG	CA	1803	1/1	0.91	0.61	-	99,99,99,99	0
54	MG	DA	3671	1/1	0.89	1.12	-	69,69,69,69	0
54	MG	BA	3553	1/1	0.94	0.17	-	72,72,72,72	0
54	MG	BA	3421	1/1	0.88	0.20	-	80,80,80,80	0
54	MG	CA	1742	1/1	0.94	0.16	-	91,91,91,91	0
54	MG	BA	2979	1/1	0.88	0.14	-	86,86,86,86	0
54	MG	CA	1745	1/1	0.89	0.19	-	77,77,77,77	0
54	MG	BA	3653	1/1	0.79	0.43	-	82,82,82,82	0
54	MG	DA	3092	1/1	0.95	0.54	-	56,56,56,56	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
54	MG	DA	3376	1/1	0.91	0.25	-	44,44,44,44	0
54	MG	AA	1838	1/1	0.85	0.64	-	77,77,77,77	0
54	MG	BA	3294	1/1	0.90	0.32	-	89,89,89,89	0
54	MG	BA	3129	1/1	0.79	0.25	-	98,98,98,98	0
54	MG	BA	3046	1/1	0.94	0.25	-	65,65,65,65	0
54	MG	BA	3363	1/1	0.93	0.33	-	113,113,113,113	0
54	MG	DA	3617	1/1	0.96	0.30	-	59,59,59,59	0
54	MG	DA	3318	1/1	0.93	0.25	-	77,77,77,77	0
54	MG	BA	3147	1/1	0.80	0.34	-	123,123,123,123	0
54	MG	BA	3357	1/1	0.76	0.38	-	80,80,80,80	0
54	MG	BA	3377	1/1	0.92	0.47	-	68,68,68,68	0
54	MG	BA	3339	1/1	0.71	0.23	-	103,103,103,103	0
54	MG	BA	3220	1/1	0.80	0.33	-	79,79,79,79	0
54	MG	AA	1877	1/1	0.89	1.48	-	102,102,102,102	0
54	MG	BA	3246	1/1	0.92	0.34	-	164,164,164,164	0
54	MG	BA	3145	1/1	0.89	0.59	-	74,74,74,74	0
54	MG	CA	1861	1/1	0.96	0.28	-	64,64,64,64	0
54	MG	CA	1842	1/1	0.82	0.41	-	103,103,103,103	0
54	MG	DA	3704	1/1	0.92	0.75	-	68,68,68,68	0
54	MG	DA	3694	1/1	0.90	0.20	-	64,64,64,64	0
54	MG	CA	1783	1/1	0.98	0.12	-	139,139,139,139	0
54	MG	DA	3241	1/1	0.85	0.09	-	76,76,76,76	0
54	MG	BA	3567	1/1	0.93	0.20	-	126,126,126,126	0
54	MG	DA	3105	1/1	0.69	0.41	-	87,87,87,87	0
54	MG	DB	235	1/1	0.81	0.41	-	91,91,91,91	0
54	MG	BA	3620	1/1	0.89	0.57	-	88,88,88,88	0
54	MG	CA	1940	1/1	0.94	0.16	-	72,72,72,72	0
54	MG	CA	1927	1/1	0.95	0.18	-	88,88,88,88	0
54	MG	DA	3676	1/1	0.80	0.23	-	92,92,92,92	0
54	MG	BA	2904	1/1	0.85	0.31	-	40,40,40,40	0
54	MG	CA	1965	1/1	0.94	0.36	-	78,78,78,78	0
54	MG	BA	3611	1/1	0.89	0.28	-	75,75,75,75	0
54	MG	DA	2970	1/1	0.96	0.25	-	29,29,29,29	0
54	MG	BA	3433	1/1	0.97	0.69	-	21,21,21,21	0
54	MG	BA	3646	1/1	0.80	0.28	-	82,82,82,82	0
54	MG	BA	3406	1/1	0.80	0.77	-	88,88,88,88	0
54	MG	DA	3209	1/1	0.74	1.22	-	94,94,94,94	0
54	MG	DA	3831	1/1	0.89	0.30	-	66,66,66,66	0
54	MG	DA	3846	1/1	0.89	0.24	-	94,94,94,94	0
54	MG	DA	3784	1/1	0.86	0.07	-	108,108,108,108	0
54	MG	BA	3075	1/1	0.74	0.13	-	60,60,60,60	0
54	MG	CD	303	1/1	0.84	0.21	-	76,76,76,76	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
54	MG	DG	202	1/1	0.63	0.13	-	102,102,102,102	0
54	MG	DA	3526	1/1	0.91	0.18	-	84,84,84,84	0
54	MG	DA	3371	1/1	0.90	0.30	-	53,53,53,53	0
54	MG	BA	3485	1/1	0.93	0.53	-	72,72,72,72	0
54	MG	DA	3046	1/1	0.85	0.35	-	37,37,37,37	0
54	MG	BA	3481	1/1	0.96	0.39	-	71,71,71,71	0
54	MG	DA	3026	1/1	0.74	0.42	-	83,83,83,83	0
54	MG	BA	3031	1/1	0.94	0.55	-	80,80,80,80	0
54	MG	CA	1771	1/1	0.90	0.17	-	98,98,98,98	0
54	MG	CA	1798	1/1	0.87	0.26	-	111,111,111,111	0
54	MG	BA	3557	1/1	0.83	0.36	-	67,67,67,67	0
54	MG	AA	1642	1/1	0.84	0.63	-	100,100,100,100	0
54	MG	AQ	201	1/1	0.72	0.30	-	90,90,90,90	0
54	MG	BA	3504	1/1	0.80	0.52	-	94,94,94,94	0
54	MG	AA	1699	1/1	0.88	0.34	-	57,57,57,57	0
54	MG	DA	3620	1/1	0.97	0.30	-	115,115,115,115	0
54	MG	DA	3260	1/1	0.86	1.20	-	90,90,90,90	0
54	MG	DA	2960	1/1	0.97	0.32	-	33,33,33,33	0
54	MG	BA	2996	1/1	0.90	0.27	-	125,125,125,125	0
54	MG	CA	1668	1/1	0.89	0.25	-	86,86,86,86	0
54	MG	DA	3758	1/1	0.91	0.31	-	60,60,60,60	0
54	MG	BA	2940	1/1	0.94	0.33	-	30,30,30,30	0
54	MG	DA	3382	1/1	0.90	0.42	-	92,92,92,92	0
54	MG	CA	1867	1/1	0.45	0.59	-	123,123,123,123	0
54	MG	CC	307	1/1	0.53	0.18	-	108,108,108,108	0
54	MG	CA	1696	1/1	0.82	0.16	-	75,75,75,75	0
54	MG	DA	3609	1/1	0.72	0.36	-	113,113,113,113	0
54	MG	BA	3394	1/1	0.92	0.53	-	95,95,95,95	0
54	MG	DA	3251	1/1	0.79	0.15	-	77,77,77,77	0
54	MG	DX	103	1/1	0.88	0.29	-	60,60,60,60	0
54	MG	BA	2927	1/1	0.86	0.24	-	25,25,25,25	0
54	MG	DA	3087	1/1	0.70	0.70	-	71,71,71,71	0
54	MG	DA	3262	1/1	0.87	0.58	-	73,73,73,73	0
54	MG	AA	1688	1/1	0.40	0.44	-	140,140,140,140	0
54	MG	DA	3086	1/1	0.61	0.46	-	71,71,71,71	0
54	MG	BA	3414	1/1	0.31	1.52	-	137,137,137,137	0
54	MG	DA	3032	1/1	0.96	0.64	-	80,80,80,80	0
54	MG	BA	2983	1/1	0.97	0.10	-	44,44,44,44	0
54	MG	DA	3585	1/1	0.81	0.24	-	154,154,154,154	0
54	MG	DA	3426	1/1	0.96	0.19	-	96,96,96,96	0
54	MG	AA	1839	1/1	0.74	0.41	-	69,69,69,69	0
54	MG	BA	3463	1/1	0.97	0.41	-	54,54,54,54	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
54	MG	BA	3388	1/1	0.87	0.52	-	66,66,66,66	0
54	MG	DA	2974	1/1	0.98	0.26	-	60,60,60,60	0
54	MG	BA	3649	1/1	0.95	0.15	-	67,67,67,67	0
54	MG	BA	3669	1/1	0.91	0.17	-	84,84,84,84	0
54	MG	DA	3475	1/1	0.62	0.68	-	120,120,120,120	0
54	MG	CA	1773	1/1	0.98	0.24	-	97,97,97,97	0
54	MG	CA	1912	1/1	0.70	0.29	-	138,138,138,138	0
54	MG	BA	3662	1/1	0.82	0.22	-	108,108,108,108	0
54	MG	BA	3518	1/1	0.95	0.69	-	81,81,81,81	0
54	MG	DA	2929	1/1	0.96	0.48	-	21,21,21,21	0
54	MG	DA	3499	1/1	0.91	0.62	-	87,87,87,87	0
54	MG	BA	3011	1/1	0.97	0.25	-	51,51,51,51	0
54	MG	CL	205	1/1	0.95	0.26	-	97,97,97,97	0
54	MG	CA	1878	1/1	0.96	0.41	-	82,82,82,82	0
54	MG	BA	3143	1/1	0.97	0.12	-	48,48,48,48	0
54	MG	DA	3823	1/1	0.96	0.51	-	84,84,84,84	0
54	MG	BA	3144	1/1	0.96	0.09	-	148,148,148,148	0
54	MG	AA	1645	1/1	0.90	0.06	-	138,138,138,138	0
54	MG	BA	3027	1/1	0.88	0.28	-	72,72,72,72	0
54	MG	CA	1700	1/1	0.94	0.16	-	129,129,129,129	0
54	MG	DA	3287	1/1	0.86	0.07	-	104,104,104,104	0
54	MG	BA	3534	1/1	0.81	0.42	-	82,82,82,82	0
54	MG	BB	202	1/1	0.85	0.44	-	48,48,48,48	0
54	MG	AA	1856	1/1	0.89	0.39	-	58,58,58,58	0
54	MG	BA	3371	1/1	0.84	0.29	-	114,114,114,114	0
54	MG	BA	3540	1/1	0.86	0.27	-	80,80,80,80	0
54	MG	BF	201	1/1	0.93	0.08	-	122,122,122,122	0
54	MG	CA	1767	1/1	0.93	0.42	-	72,72,72,72	0
54	MG	DA	3791	1/1	0.88	0.28	-	108,108,108,108	0
54	MG	DA	3640	1/1	0.98	0.25	-	24,24,24,24	0
54	MG	DA	3309	1/1	0.93	0.21	-	127,127,127,127	0
54	MG	CA	1886	1/1	0.70	0.32	-	105,105,105,105	0
54	MG	AA	1767	1/1	0.90	0.14	-	118,118,118,118	0
54	MG	DA	3509	1/1	0.81	0.14	-	148,148,148,148	0
54	MG	DA	3159	1/1	0.93	0.43	-	106,106,106,106	0
54	MG	DA	2966	1/1	0.97	0.19	-	39,39,39,39	0
54	MG	BA	3234	1/1	0.86	0.42	-	73,73,73,73	0
54	MG	BA	3656	1/1	0.93	0.22	-	107,107,107,107	0
54	MG	DA	2911	1/1	0.94	0.38	-	11,11,11,11	0
54	MG	DA	3143	1/1	0.83	0.25	-	67,67,67,67	0
54	MG	BA	3128	1/1	0.98	0.18	-	65,65,65,65	0
54	MG	AA	1845	1/1	0.96	0.06	-	71,71,71,71	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
54	MG	BA	3321	1/1	0.98	0.22	-	179,179,179,179	0
54	MG	CA	1633	1/1	0.93	0.21	-	85,85,85,85	0
54	MG	BA	3375	1/1	0.92	0.28	-	62,62,62,62	0
54	MG	BA	3462	1/1	0.79	0.61	-	80,80,80,80	0
54	MG	DA	2991	1/1	0.99	0.21	-	36,36,36,36	0
54	MG	DA	2930	1/1	0.99	0.49	-	34,34,34,34	0
54	MG	DA	3471	1/1	0.56	0.30	-	104,104,104,104	0
54	MG	DA	3648	1/1	0.80	0.48	-	60,60,60,60	0
54	MG	DS	202	1/1	0.95	0.52	-	62,62,62,62	0
54	MG	BA	3604	1/1	0.69	0.29	-	89,89,89,89	0
54	MG	AA	1710	1/1	0.87	0.52	-	84,84,84,84	0
54	MG	AA	1829	1/1	0.90	0.13	-	64,64,64,64	0
54	MG	DA	3774	1/1	0.94	0.30	-	74,74,74,74	0
54	MG	AA	1696	1/1	0.94	0.05	-	120,120,120,120	0
54	MG	DA	3451	1/1	0.93	0.35	-	80,80,80,80	0
54	MG	BA	3642	1/1	0.83	0.18	-	98,98,98,98	0
54	MG	DA	3243	1/1	0.98	0.12	-	59,59,59,59	0
54	MG	DA	3391	1/1	0.94	0.37	-	82,82,82,82	0
54	MG	BA	3498	1/1	0.81	0.97	-	67,67,67,67	0
54	MG	BA	2994	1/1	0.84	0.55	-	90,90,90,90	0
54	MG	CA	1966	1/1	0.92	0.30	-	97,97,97,97	0
54	MG	DB	239	1/1	0.73	0.17	-	107,107,107,107	0
54	MG	CA	1646	1/1	0.95	0.21	-	115,115,115,115	0
54	MG	DA	3367	1/1	0.74	0.92	-	97,97,97,97	0
54	MG	DA	3517	1/1	0.78	0.21	-	102,102,102,102	0
54	MG	AA	1643	1/1	0.86	0.19	-	63,63,63,63	0
54	MG	BA	3618	1/1	0.90	0.22	-	76,76,76,76	0
54	MG	DA	3088	1/1	0.90	0.09	-	83,83,83,83	0
54	MG	DA	3481	1/1	0.97	0.29	-	77,77,77,77	0
54	MG	CA	1615	1/1	0.94	0.20	-	70,70,70,70	0
54	MG	BA	3648	1/1	0.82	0.31	-	63,63,63,63	0
54	MG	BA	3171	1/1	0.90	0.10	-	127,127,127,127	0
54	MG	BA	3533	1/1	0.93	0.20	-	83,83,83,83	0
54	MG	DA	3464	1/1	0.88	0.12	-	133,133,133,133	0
54	MG	DA	3007	1/1	0.88	0.36	-	44,44,44,44	0
54	MG	DA	3017	1/1	0.96	0.33	-	38,38,38,38	0
54	MG	CA	1736	1/1	0.82	0.25	-	73,73,73,73	0
54	MG	DA	3463	1/1	0.71	0.42	-	69,69,69,69	0
54	MG	BA	3351	1/1	0.90	0.76	-	116,116,116,116	0
54	MG	DA	2909	1/1	0.95	0.61	-	22,22,22,22	0
54	MG	CA	1740	1/1	0.76	0.14	-	117,117,117,117	0
54	MG	BA	3663	1/1	0.90	0.31	-	89,89,89,89	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
54	MG	BV	203	1/1	0.92	0.09	-	104,104,104,104	0
54	MG	DA	3291	1/1	0.82	0.27	-	57,57,57,57	0
54	MG	DA	3424	1/1	0.76	0.19	-	90,90,90,90	0
54	MG	BA	3520	1/1	0.63	0.87	-	95,95,95,95	0
54	MG	BA	3578	1/1	0.93	0.10	-	74,74,74,74	0
54	MG	BA	3621	1/1	0.83	0.58	-	107,107,107,107	0
54	MG	BA	2997	1/1	0.94	0.71	-	81,81,81,81	0
54	MG	BA	3526	1/1	0.89	0.18	-	66,66,66,66	0
54	MG	DA	3421	1/1	0.89	0.15	-	63,63,63,63	0
54	MG	AA	1831	1/1	0.95	0.40	-	67,67,67,67	0
54	MG	CA	1759	1/1	0.90	0.32	-	89,89,89,89	0
54	MG	DA	3296	1/1	0.96	0.24	-	70,70,70,70	0
54	MG	BA	3309	1/1	0.90	0.26	-	84,84,84,84	0
54	MG	BA	3660	1/1	0.95	0.15	-	119,119,119,119	0
54	MG	BA	2941	1/1	0.96	0.34	-	31,31,31,31	0
54	MG	CA	1775	1/1	0.94	0.17	-	64,64,64,64	0
54	MG	DA	3781	1/1	0.80	0.34	-	55,55,55,55	0
54	MG	AA	1807	1/1	0.82	0.51	-	76,76,76,76	0
54	MG	DA	3772	1/1	0.98	0.21	-	114,114,114,114	0
54	MG	BA	3163	1/1	0.79	0.45	-	87,87,87,87	0
54	MG	DA	3647	1/1	0.99	0.75	-	54,54,54,54	0
54	MG	CA	1889	1/1	0.67	0.33	-	74,74,74,74	0
54	MG	CA	1681	1/1	0.97	0.20	-	66,66,66,66	0
54	MG	AA	1761	1/1	0.96	0.10	-	68,68,68,68	0
54	MG	BA	3457	1/1	0.94	0.37	-	48,48,48,48	0
54	MG	CA	1644	1/1	0.91	0.15	-	94,94,94,94	0
54	MG	AA	1705	1/1	0.70	0.57	-	62,62,62,62	0
54	MG	BA	3679	1/1	0.96	0.29	-	75,75,75,75	0
54	MG	DA	3284	1/1	0.96	0.20	-	79,79,79,79	0
54	MG	DA	3803	1/1	0.56	0.51	-	123,123,123,123	0
54	MG	BA	3316	1/1	0.87	0.44	-	97,97,97,97	0
54	MG	DA	3029	1/1	0.80	0.37	-	74,74,74,74	0
54	MG	BA	3402	1/1	0.75	0.33	-	90,90,90,90	0
54	MG	DA	3297	1/1	0.88	0.55	-	76,76,76,76	0
54	MG	BA	3360	1/1	0.95	0.20	-	75,75,75,75	0
54	MG	DA	3575	1/1	0.95	0.38	-	85,85,85,85	0
54	MG	DA	3443	1/1	0.97	0.15	-	92,92,92,92	0
54	MG	BB	203	1/1	0.89	0.75	-	74,74,74,74	0
54	MG	AA	1832	1/1	0.96	0.46	-	66,66,66,66	0
54	MG	BA	3381	1/1	0.90	0.63	-	82,82,82,82	0
54	MG	DA	3654	1/1	0.97	0.28	-	59,59,59,59	0
54	MG	DA	3051	1/1	0.92	0.30	-	60,60,60,60	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
54	MG	AV	6201	1/1	0.91	0.57	-	67,67,67,67	0
54	MG	BA	3259	1/1	0.92	0.08	-	105,105,105,105	0
54	MG	DA	3627	1/1	0.88	0.60	-	63,63,63,63	0
54	MG	CA	1866	1/1	0.96	0.21	-	102,102,102,102	0
54	MG	BA	3671	1/1	0.93	0.25	-	100,100,100,100	0
54	MG	CA	1879	1/1	0.95	0.17	-	76,76,76,76	0
54	MG	BA	3563	1/1	0.74	0.38	-	58,58,58,58	0
54	MG	BA	3474	1/1	0.94	0.33	-	72,72,72,72	0
54	MG	BA	3569	1/1	0.80	0.13	-	116,116,116,116	0
54	MG	BA	3295	1/1	0.89	0.30	-	60,60,60,60	0
54	MG	DA	3127	1/1	0.85	0.45	-	76,76,76,76	0
54	MG	DA	3541	1/1	0.92	0.19	-	76,76,76,76	0
54	MG	DA	3109	1/1	0.74	0.32	-	88,88,88,88	0
54	MG	BA	3313	1/1	0.94	0.18	-	67,67,67,67	0
54	MG	BA	3092	1/1	0.82	0.70	-	123,123,123,123	0
54	MG	DA	3073	1/1	0.96	0.24	-	69,69,69,69	0
54	MG	DA	3832	1/1	0.73	0.30	-	77,77,77,77	0
54	MG	BA	3087	1/1	0.92	1.47	-	84,84,84,84	0
54	MG	DA	3390	1/1	0.76	0.78	-	72,72,72,72	0
54	MG	BA	3666	1/1	0.93	0.19	-	72,72,72,72	0
54	MG	DA	3366	1/1	0.80	0.39	-	73,73,73,73	0
54	MG	BA	3449	1/1	0.94	0.57	-	31,31,31,31	0
54	MG	AA	1830	1/1	0.08	0.86	-	105,105,105,105	0
54	MG	BA	3458	1/1	0.97	0.11	-	77,77,77,77	0
54	MG	DA	3746	1/1	0.95	0.11	-	94,94,94,94	0
54	MG	BA	3509	1/1	0.92	0.17	-	64,64,64,64	0
54	MG	DA	3630	1/1	0.92	0.23	-	101,101,101,101	0
54	MG	BA	3176	1/1	0.92	0.29	-	53,53,53,53	0
54	MG	BA	3276	1/1	0.97	0.07	-	67,67,67,67	0
54	MG	AA	1844	1/1	0.37	1.02	-	121,121,121,121	0
54	MG	CA	1658	1/1	0.88	0.26	-	94,94,94,94	0
54	MG	BN	201	1/1	0.96	0.28	-	49,49,49,49	0
54	MG	DA	3729	1/1	0.93	0.17	-	81,81,81,81	0
54	MG	CA	1796	1/1	0.81	0.38	-	87,87,87,87	0
54	MG	AA	1857	1/1	0.88	0.14	-	103,103,103,103	0
54	MG	DA	3270	1/1	0.93	0.24	-	82,82,82,82	0
54	MG	BA	2972	1/1	0.85	0.33	-	61,61,61,61	0
54	MG	DA	3468	1/1	0.65	0.49	-	79,79,79,79	0
54	MG	DA	3047	1/1	0.78	0.46	-	74,74,74,74	0
54	MG	DA	3745	1/1	0.96	0.25	-	109,109,109,109	0
54	MG	BA	3010	1/1	0.94	0.07	-	73,73,73,73	0
54	MG	BB	208	1/1	0.96	0.11	-	67,67,67,67	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
54	MG	BA	3487	1/1	0.84	0.41	-	90,90,90,90	0
54	MG	BA	3588	1/1	0.82	0.74	-	52,52,52,52	0
54	MG	BA	3262	1/1	0.83	0.14	-	75,75,75,75	0
54	MG	AA	1639	1/1	0.91	0.60	-	54,54,54,54	0
54	MG	CA	1638	1/1	0.84	0.06	-	101,101,101,101	0
54	MG	DA	3519	1/1	0.96	0.38	-	74,74,74,74	0
54	MG	CA	1810	1/1	0.92	0.09	-	99,99,99,99	0
54	MG	DA	3527	1/1	0.94	0.17	-	104,104,104,104	0
54	MG	DA	3656	1/1	0.96	0.29	-	36,36,36,36	0
54	MG	BA	3034	1/1	0.98	0.98	-	69,69,69,69	0
54	MG	BA	3191	1/1	0.94	0.95	-	90,90,90,90	0
54	MG	DA	2998	1/1	0.95	0.26	-	36,36,36,36	0
54	MG	CA	1967	1/1	0.83	0.33	-	93,93,93,93	0
54	MG	DA	3372	1/1	0.98	0.16	-	90,90,90,90	0
54	MG	BA	3332	1/1	0.91	0.20	-	65,65,65,65	0
54	MG	DA	3379	1/1	0.95	0.04	-	78,78,78,78	0
54	MG	BA	3672	1/1	0.92	0.30	-	92,92,92,92	0
54	MG	BA	3174	1/1	0.86	0.30	-	67,67,67,67	0
54	MG	AA	1612	1/1	0.92	0.58	-	39,39,39,39	0
54	MG	DA	3770	1/1	0.85	0.17	-	79,79,79,79	0
54	MG	BA	3202	1/1	0.79	0.34	-	126,126,126,126	0
54	MG	DA	3754	1/1	0.30	0.37	-	121,121,121,121	0
54	MG	AA	1862	1/1	0.87	0.66	-	84,84,84,84	0
54	MG	BA	3435	1/1	0.98	0.38	-	29,29,29,29	0
54	MG	BA	3641	1/1	0.93	0.15	-	95,95,95,95	0
54	MG	BA	3311	1/1	0.98	0.17	-	87,87,87,87	0
54	MG	DA	3062	1/1	0.92	0.28	-	129,129,129,129	0
54	MG	BA	3288	1/1	0.77	0.53	-	113,113,113,113	0
54	MG	AA	1825	1/1	0.87	0.98	-	79,79,79,79	0
54	MG	DA	3696	1/1	0.92	0.15	-	82,82,82,82	0
54	MG	BA	3281	1/1	0.96	0.72	-	123,123,123,123	0
54	MG	DN	201	1/1	0.95	0.07	-	88,88,88,88	0
54	MG	BA	2958	1/1	0.91	0.60	-	46,46,46,46	0
54	MG	DA	3452	1/1	0.77	0.53	-	74,74,74,74	0
54	MG	CA	1791	1/1	0.92	0.84	-	93,93,93,93	0
54	MG	D3	101	1/1	0.85	0.14	-	90,90,90,90	0
54	MG	DA	3677	1/1	0.93	0.23	-	60,60,60,60	0
54	MG	DA	3459	1/1	0.99	0.16	-	61,61,61,61	0
54	MG	BA	3236	1/1	0.91	0.23	-	56,56,56,56	0
54	MG	DA	3544	1/1	0.94	0.27	-	53,53,53,53	0
54	MG	CF	202	1/1	0.82	0.22	-	83,83,83,83	0
54	MG	DA	3427	1/1	0.88	0.38	-	85,85,85,85	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
54	MG	CA	1627	1/1	0.77	1.09	-	99,99,99,99	0
54	MG	BU	202	1/1	0.78	0.25	-	74,74,74,74	0
54	MG	BA	3193	1/1	0.98	0.29	-	81,81,81,81	0
54	MG	DA	3610	1/1	0.89	0.57	-	81,81,81,81	0
54	MG	CA	1845	1/1	0.75	1.93	-	100,100,100,100	0
54	MG	DA	2950	1/1	0.98	0.23	-	33,33,33,33	0
54	MG	CL	206	1/1	0.77	0.16	-	87,87,87,87	0
54	MG	CA	1838	1/1	0.97	0.14	-	94,94,94,94	0
54	MG	BA	3221	1/1	0.88	0.45	-	141,141,141,141	0
54	MG	CA	1920	1/1	0.98	0.28	-	87,87,87,87	0
54	MG	CA	1626	1/1	0.88	0.54	-	68,68,68,68	0
54	MG	DA	3793	1/1	0.85	0.25	-	106,106,106,106	0
54	MG	AA	1700	1/1	0.59	0.38	-	114,114,114,114	0
54	MG	CA	1753	1/1	0.91	0.32	-	97,97,97,97	0
54	MG	CA	1852	1/1	0.82	0.38	-	101,101,101,101	0
54	MG	DA	3717	1/1	0.91	0.26	-	65,65,65,65	0
54	MG	BA	3223	1/1	0.94	0.16	-	76,76,76,76	0
54	MG	DA	3722	1/1	0.97	0.17	-	138,138,138,138	0
54	MG	CV	6206	1/1	0.93	0.08	-	107,107,107,107	0
54	MG	DA	2987	1/1	0.91	0.27	-	28,28,28,28	0
54	MG	AA	1806	1/1	0.89	0.34	-	65,65,65,65	0
54	MG	DB	210	1/1	0.94	0.10	-	92,92,92,92	0
54	MG	DA	3316	1/1	0.93	0.16	-	94,94,94,94	0
54	MG	BA	2921	1/1	0.98	0.27	-	17,17,17,17	0
54	MG	BA	2985	1/1	0.89	0.54	-	43,43,43,43	0
54	MG	CA	1719	1/1	0.81	0.43	-	78,78,78,78	0
54	MG	CA	1871	1/1	0.97	0.49	-	76,76,76,76	0
54	MG	BA	3454	1/1	0.93	0.19	-	41,41,41,41	0
54	MG	DA	3491	1/1	0.82	1.22	-	99,99,99,99	0
54	MG	DB	232	1/1	0.85	0.21	-	84,84,84,84	0
54	MG	CA	1818	1/1	0.82	0.42	-	91,91,91,91	0
54	MG	AA	1678	1/1	0.90	0.20	-	78,78,78,78	0
54	MG	CQ	201	1/1	0.84	0.38	-	103,103,103,103	0
54	MG	DA	3785	1/1	0.89	0.25	-	109,109,109,109	0
54	MG	AA	1874	1/1	0.81	0.11	-	87,87,87,87	0
54	MG	BA	3670	1/1	0.83	0.73	-	90,90,90,90	0
54	MG	BA	2938	1/1	0.97	0.46	-	51,51,51,51	0
54	MG	DA	3110	1/1	0.95	0.24	-	65,65,65,65	0
54	MG	BA	3102	1/1	0.88	0.35	-	72,72,72,72	0
54	MG	AE	203	1/1	0.78	0.12	-	91,91,91,91	0
54	MG	BA	3146	1/1	0.87	0.05	-	99,99,99,99	0
54	MG	DA	3039	1/1	0.95	0.11	-	70,70,70,70	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
54	MG	DA	3730	1/1	0.64	0.28	-	61,61,61,61	0
54	MG	DA	3457	1/1	0.85	0.28	-	86,86,86,86	0
54	MG	DA	3266	1/1	0.93	0.16	-	85,85,85,85	0
54	MG	BA	3042	1/1	0.76	0.88	-	92,92,92,92	0
54	MG	DA	3128	1/1	0.91	0.46	-	85,85,85,85	0
54	MG	DA	3053	1/1	0.97	0.47	-	63,63,63,63	0
54	MG	DA	3432	1/1	0.63	0.27	-	84,84,84,84	0
54	MG	CA	1859	1/1	0.88	0.25	-	93,93,93,93	0
54	MG	BA	3149	1/1	0.92	0.36	-	65,65,65,65	0
54	MG	CA	1634	1/1	0.91	0.42	-	78,78,78,78	0
54	MG	BA	3074	1/1	0.97	0.38	-	87,87,87,87	0
54	MG	DA	3381	1/1	0.95	0.23	-	49,49,49,49	0
54	MG	AA	1666	1/1	0.74	0.56	-	93,93,93,93	0
54	MG	CA	1789	1/1	0.93	0.21	-	103,103,103,103	0
54	MG	BA	3273	1/1	0.78	0.36	-	73,73,73,73	0
54	MG	DA	3756	1/1	0.72	0.68	-	80,80,80,80	0
54	MG	CA	1727	1/1	0.91	0.15	-	76,76,76,76	0
54	MG	DA	3844	1/1	0.70	0.21	-	132,132,132,132	0
54	MG	DF	202	1/1	0.97	0.19	-	129,129,129,129	0
54	MG	BI	101	1/1	0.59	0.33	-	109,109,109,109	0
54	MG	AV	6204	1/1	0.54	0.13	-	124,124,124,124	0
54	MG	BT	102	1/1	0.90	0.23	-	66,66,66,66	0
54	MG	DA	3192	1/1	0.94	0.18	-	49,49,49,49	0
54	MG	DA	3489	1/1	0.93	0.29	-	97,97,97,97	0
54	MG	BA	2901	1/1	0.99	0.63	-	8,8,8,8	0
54	MG	CA	1916	1/1	0.88	0.38	-	118,118,118,118	0
54	MG	DA	3215	1/1	0.84	0.39	-	73,73,73,73	0
54	MG	DA	3205	1/1	0.71	0.51	-	97,97,97,97	0
54	MG	AA	1734	1/1	0.76	0.21	-	96,96,96,96	0
54	MG	AA	1722	1/1	0.82	0.51	-	53,53,53,53	0
54	MG	DA	3141	1/1	0.95	0.44	-	57,57,57,57	0
54	MG	CV	6204	1/1	0.93	0.13	-	112,112,112,112	0
54	MG	DA	3685	1/1	0.82	0.39	-	82,82,82,82	0
54	MG	BA	3605	1/1	0.76	0.42	-	69,69,69,69	0
54	MG	DA	3041	1/1	0.91	0.23	-	76,76,76,76	0
54	MG	BA	3523	1/1	0.89	1.59	-	99,99,99,99	0
54	MG	D4	103	1/1	0.95	0.27	-	73,73,73,73	0
54	MG	DA	3170	1/1	0.95	0.41	-	56,56,56,56	0
54	MG	CA	1953	1/1	0.93	0.24	-	120,120,120,120	0
54	MG	CA	1846	1/1	0.93	0.30	-	135,135,135,135	0
54	MG	DA	3852	1/1	0.63	0.20	-	80,80,80,80	0
54	MG	BA	3584	1/1	0.96	0.15	-	83,83,83,83	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
54	MG	DA	3618	1/1	0.76	0.54	-	129,129,129,129	0
54	MG	DA	3419	1/1	0.94	0.13	-	60,60,60,60	0
54	MG	AA	1698	1/1	0.99	0.13	-	88,88,88,88	0
54	MG	DA	3233	1/1	0.92	0.29	-	67,67,67,67	0
54	MG	BA	3379	1/1	0.96	0.21	-	86,86,86,86	0
54	MG	DA	3639	1/1	0.97	0.35	-	26,26,26,26	0
54	MG	BA	3310	1/1	0.75	0.44	-	109,109,109,109	0
54	MG	DA	3394	1/1	0.96	0.53	-	110,110,110,110	0
54	MG	BA	3418	1/1	0.52	0.42	-	88,88,88,88	0
54	MG	BA	3658	1/1	0.96	0.29	-	68,68,68,68	0
54	MG	CA	1782	1/1	0.92	0.07	-	80,80,80,80	0
54	MG	CA	1761	1/1	0.94	0.95	-	88,88,88,88	0
54	MG	AA	1787	1/1	0.67	0.51	-	68,68,68,68	0
54	MG	BA	3506	1/1	0.95	0.36	-	89,89,89,89	0
54	MG	DA	3416	1/1	0.94	0.25	-	101,101,101,101	0
54	MG	DA	3118	1/1	0.89	0.43	-	50,50,50,50	0
54	MG	BA	3308	1/1	0.70	0.21	-	99,99,99,99	0
54	MG	CA	1751	1/1	0.87	0.09	-	117,117,117,117	0
54	MG	DA	3850	1/1	0.87	0.41	-	52,52,52,52	0
54	MG	DA	2936	1/1	0.90	0.12	-	42,42,42,42	0
54	MG	AA	1813	1/1	0.96	0.12	-	76,76,76,76	0
54	MG	BA	3006	1/1	0.87	0.57	-	61,61,61,61	0
54	MG	CA	1741	1/1	0.93	0.28	-	100,100,100,100	0
54	MG	AA	1843	1/1	0.67	1.09	-	126,126,126,126	0
54	MG	AA	1759	1/1	0.82	0.48	-	110,110,110,110	0
54	MG	BA	3650	1/1	0.89	0.31	-	69,69,69,69	0
54	MG	BA	3033	1/1	0.75	0.43	-	77,77,77,77	0
54	MG	BA	3121	1/1	0.91	0.22	-	69,69,69,69	0
54	MG	DA	3428	1/1	0.80	0.38	-	71,71,71,71	0
54	MG	BA	2964	1/1	0.94	0.35	-	43,43,43,43	0
54	MG	DA	3111	1/1	0.89	0.13	-	84,84,84,84	0
54	MG	AA	1668	1/1	0.93	0.17	-	77,77,77,77	0
54	MG	BA	3157	1/1	0.87	0.18	-	94,94,94,94	0
54	MG	BA	3069	1/1	0.95	0.26	-	52,52,52,52	0
54	MG	AA	1814	1/1	0.89	0.12	-	82,82,82,82	0
54	MG	DA	3036	1/1	0.90	0.21	-	72,72,72,72	0
54	MG	BA	3348	1/1	0.89	0.38	-	107,107,107,107	0
54	MG	DA	3810	1/1	0.97	0.05	-	102,102,102,102	0
54	MG	BA	3573	1/1	0.94	0.17	-	58,58,58,58	0
54	MG	DA	3473	1/1	0.94	0.47	-	61,61,61,61	0
54	MG	BY	102	1/1	0.97	0.11	-	110,110,110,110	0
54	MG	DA	3441	1/1	0.93	0.15	-	60,60,60,60	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
54	MG	CA	1679	1/1	0.88	0.53	-	89,89,89,89	0
54	MG	BA	3109	1/1	0.83	0.48	-	77,77,77,77	0
54	MG	AL	201	1/1	0.92	0.21	-	109,109,109,109	0
54	MG	DA	3090	1/1	0.89	0.42	-	59,59,59,59	0
54	MG	BA	3035	1/1	0.97	0.23	-	57,57,57,57	0
54	MG	CA	1865	1/1	0.98	0.57	-	93,93,93,93	0
54	MG	CA	1968	1/1	0.90	0.15	-	87,87,87,87	0
54	MG	DA	2922	1/1	0.90	0.46	-	37,37,37,37	0
54	MG	CV	6203	1/1	0.85	0.38	-	80,80,80,80	0
54	MG	BA	3214	1/1	0.88	0.80	-	79,79,79,79	0
54	MG	BA	3503	1/1	0.89	0.83	-	90,90,90,90	0
54	MG	DA	2988	1/1	0.80	0.43	-	23,23,23,23	0
54	MG	DA	3335	1/1	0.94	0.10	-	46,46,46,46	0
54	MG	BA	3536	1/1	0.82	0.34	-	83,83,83,83	0
54	MG	CA	1870	1/1	0.97	0.20	-	94,94,94,94	0
54	MG	AA	1683	1/1	0.97	0.46	-	64,64,64,64	0
54	MG	BA	3272	1/1	0.84	0.79	-	81,81,81,81	0
54	MG	CA	1686	1/1	0.96	0.07	-	82,82,82,82	0
54	MG	BA	3210	1/1	0.91	0.06	-	86,86,86,86	0
54	MG	DA	3389	1/1	0.90	0.17	-	76,76,76,76	0
54	MG	CO	101	1/1	0.38	0.19	-	127,127,127,127	0
54	MG	AC	301	1/1	0.83	0.22	-	104,104,104,104	0
54	MG	BA	3448	1/1	0.89	0.47	-	51,51,51,51	0
54	MG	CA	1695	1/1	0.96	0.18	-	65,65,65,65	0
54	MG	DA	3091	1/1	0.91	0.32	-	127,127,127,127	0
54	MG	BA	3101	1/1	0.81	0.88	-	74,74,74,74	0
54	MG	AA	1827	1/1	0.92	0.28	-	71,71,71,71	0
54	MG	AA	1822	1/1	0.94	0.21	-	92,92,92,92	0
54	MG	DA	3429	1/1	0.97	0.15	-	79,79,79,79	0
54	MG	CA	1704	1/1	0.13	0.49	-	140,140,140,140	0
54	MG	AA	1842	1/1	0.93	0.89	-	69,69,69,69	0
54	MG	DA	3662	1/1	0.65	0.80	-	71,71,71,71	0
54	MG	BA	3631	1/1	0.88	0.12	-	85,85,85,85	0
54	MG	DA	3248	1/1	0.72	0.16	-	80,80,80,80	0
54	MG	CA	1893	1/1	0.73	0.15	-	127,127,127,127	0

## 6.5 Other polymers

There are no such residues in this entry.