



wwPDB EM Map/Model Validation Report ⓘ

Apr 10, 2016 – 03:15 PM BST

PDB ID : 4V7H
EMDB ID: : EMD-1345
Title : Structure of the 80S rRNA and proteins and P/E tRNA for eukaryotic ribosome based on cryo-EM map of *Thermomyces lanuginosus* ribosome at 8.9Å resolution
Authors : Taylor, D.J.; Devkota, B.; Huang, A.D.; Topf, M.; Narayanan, E.; Sali, A.; Harvey, S.C.; Frank, J.
Deposited on : 2009-09-22
Resolution : 8.90 Å(reported)

This is a wwPDB EM Map/Model Validation Report for a publicly released PDB/EMDB entry.
For rigid body fitted models, validation errors reported here could stem from errors in the original structure(s) used in the fitting.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<http://wwpdb.org/validation/2016/EMValidationReportHelp>

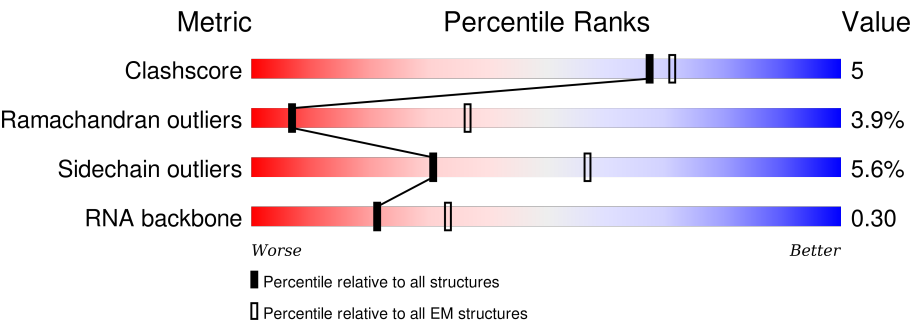
MolProbity : 4.02b-467
Mogul : 1.7.1 (RC1), CSD as537be (2016)
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : trunk27241

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
ELECTRON MICROSCOPY

The reported resolution of this entry is 8.90 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)	EM structures (#Entries)
Clashscore	114402	924
Ramachandran outliers	111179	726
Sidechain outliers	111093	686
RNA backbone	3027	244

The table below summarises the geometric issues observed across the polymeric chains. The red, orange, yellow and green segments on the bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$

Mol	Chain	Length	Quality of chain
1	AA	1761	<div><div>72%</div><div>24%</div><div></div></div>
2	AB	193	<div><div>97%</div><div></div><div></div></div>
3	AC	188	<div><div>93%</div><div>7%</div><div></div></div>
4	AD	158	<div><div>65%</div><div>12%</div><div>22%</div></div>
5	AE	162	<div><div>93%</div><div>6%</div><div></div></div>
6	AG	186	<div><div>91%</div><div>8%</div><div></div></div>
7	AH	125	<div><div>89%</div><div>7%</div><div></div></div>
8	AI	138	<div><div>83%</div><div>11%</div><div></div></div>


















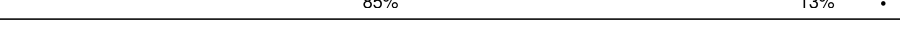


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Mol	Chain	Length	Quality of chain
9	AJ	96	
10	AK	125	
11	AL	118	
12	AM	130	
13	AN	50	
14	AO	84	
15	AQ	80	
16	AS	71	
17	AR	313	
18	AT	141	
19	A7	76	
20	B0	109	
21	B1	48	
22	B2	98	
23	B8	118	
24	B9	72	
25	BA	213	
26	BB	243	
27	BC	362	
28	BD	257	
29	BE	237	
30	BF	213	
31	BG	113	
32	BH	179	
33	BI	165	

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Mol	Chain	Length	Quality of chain
34	BJ	151	 89% 10% .
35	BK	138	 89% 10% .
36	BL	192	 90% 9% .
37	BM	178	 77% 19% . .
38	BN	150	 93% 7%
39	BO	121	 94% 5% .
40	BP	176	 83% 16% .
41	BQ	116	 91% 9%
42	BR	131	 86% 13% .
43	BS	45	 82% 16% .
44	BT	80	 86% 13% .
45	BU	116	 95% . .
46	BV	142	 83% 15% .
47	BW	79	 91% 6% .
48	BX	86	 86% 14%
49	BY	52	 90% 8% .
50	BZ	92	 85% 13% .
51	B3	113	 . 60% 36%
52	B4	157	 . 60% 36% .
53	B5	3170	 . 65% 31% .

2 Entry composition

There are 53 unique types of molecules in this entry. The entry contains 165754 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, 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 18S rRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	AA	1761	Total	C	N	O	P	0	3
			37458	16745	6626	12327	1760		

- Molecule 2 is a protein called 40S ribosomal protein S0(A).

Mol	Chain	Residues	Atoms					AltConf	Trace
2	AB	193	Total	C	N	O	S	0	0
			1500	958	269	271	2		

- Molecule 3 is a protein called 40S ribosomal protein S3.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	AC	188	Total	C	N	O	S	0	0
			1469	929	271	263	6		

- Molecule 4 is a protein called 40S ribosomal protein S9(A).

Mol	Chain	Residues	Atoms					AltConf	Trace
4	AD	124	Total	C	N	O	S	0	0
			1018	647	189	181	1		

- Molecule 5 is a protein called 40S ribosomal protein S2.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	AE	162	Total	C	N	O	S	0	0
			1207	765	222	218	2		

- Molecule 6 is a protein called 40S ribosomal protein S5.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	AG	186	Total	C	N	O	S	0	0
			1456	908	277	268	3		

- Molecule 7 is a protein called 40S ribosomal protein S22(A).

Mol	Chain	Residues	Atoms					AltConf	Trace
7	AH	125	Total	C	N	O	S	0	0
			992	634	181	174	3		

- Molecule 8 is a protein called 40S ribosomal protein S16(A).

Mol	Chain	Residues	Atoms					AltConf	Trace
8	AI	138	Total	C	N	O	S	0	0
			1087	695	200	192			

- Molecule 9 is a protein called 40S ribosomal protein S20.

Mol	Chain	Residues	Atoms					AltConf	Trace
9	AJ	96	Total	C	N	O	S	0	0
			771	487	140	143	1		

- Molecule 10 is a protein called 40S ribosomal protein S14(A).

Mol	Chain	Residues	Atoms					AltConf	Trace
10	AK	125	Total	C	N	O	S	0	0
			924	566	179	176	3		

- Molecule 11 is a protein called 40S ribosomal protein S23(A).

Mol	Chain	Residues	Atoms					AltConf	Trace
11	AL	118	Total	C	N	O	S	0	0
			906	579	166	159	2		

- Molecule 12 is a protein called 40S ribosomal protein S18.

Mol	Chain	Residues	Atoms					AltConf	Trace
12	AM	130	Total	C	N	O	S	0	0
			1077	669	217	189	2		

- Molecule 13 is a protein called 40S ribosomal protein S29(A).

Mol	Chain	Residues	Atoms					AltConf	Trace
13	AN	50	Total	C	N	O	S	0	0
			417	258	87	68	4		

- Molecule 14 is a protein called 40S ribosomal protein S13.

Mol	Chain	Residues	Atoms					AltConf	Trace
14	AO	84	Total	C	N	O	S	0	0
			694	446	129	118	1		

- Molecule 15 is a protein called 40S ribosomal protein S11.

Mol	Chain	Residues	Atoms					AltConf	Trace
15	AQ	80	Total	C	N	O	S	0	0
			643	410	127	104	2		

- Molecule 16 is a protein called 40S ribosomal protein S15.

Mol	Chain	Residues	Atoms					AltConf	Trace
16	AS	71	Total	C	N	O	S	0	0
			548	348	101	93	6		

- Molecule 17 is a protein called RACK1 protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
17	AR	313	Total	C	N	O	S	0	0
			2410	1526	413	463	8		

- Molecule 18 is a protein called s19e protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
18	AT	141	Total	C	N	O	S	0	0
			1102	687	206	207	2		

- Molecule 19 is a RNA chain called P/E tRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
19	A7	76	Total	C	N	O	P	0	0
			1648	746	294	533	75		

- Molecule 20 is a protein called 60S ribosomal protein L32.

Mol	Chain	Residues	Atoms					AltConf	Trace
20	B0	109	Total	C	N	O	S	0	0
			881	555	176	149	1		

- Molecule 21 is a protein called 60S ribosomal protein L39.

Mol	Chain	Residues	Atoms					AltConf	Trace
21	B1	48	Total	C	N	O	S	0	0
			424	263	95	64	2		

- Molecule 22 is a protein called 60S ribosomal protein L30e.

Mol	Chain	Residues	Atoms					AltConf	Trace
22	B2	98	Total	C	N	O	S	0	0
			752	484	125	142	1		

- Molecule 23 is a protein called 60S ribosomal protein LP0.

Mol	Chain	Residues	Atoms					AltConf	Trace
23	B8	118	Total	C	N	O	S	0	0
			947	609	167	168	3		

- Molecule 24 is a protein called 60S ribosomal protein L43.

Mol	Chain	Residues	Atoms					AltConf	Trace
24	B9	72	Total	C	N	O	S	0	0
			539	332	104	98	5		

- Molecule 25 is a protein called 60S ribosomal protein L1.

Mol	Chain	Residues	Atoms					AltConf	Trace
25	BA	213	Total	C	N	O	S	0	0
			1683	1074	294	306	9		

- Molecule 26 is a protein called 60S ribosomal protein L2.

Mol	Chain	Residues	Atoms					AltConf	Trace
26	BB	243	Total	C	N	O	S	0	0
			1848	1150	374	323	1		

- Molecule 27 is a protein called 60S ribosomal protein L3.

Mol	Chain	Residues	Atoms					AltConf	Trace
27	BC	362	Total	C	N	O	S	0	0
			2887	1833	545	502	7		

- Molecule 28 is a protein called 60S ribosomal protein L4(B).

Mol	Chain	Residues	Atoms					AltConf	Trace
28	BD	257	Total	C	N	O	S	0	0
			1950	1226	375	346	3		

- Molecule 29 is a protein called 60S ribosomal protein L5.

Mol	Chain	Residues	Atoms					AltConf	Trace
29	BE	237	Total	C	N	O	S	0	0
			1913	1210	329	372	2		

- Molecule 30 is a protein called 60S ribosomal protein L7(A).

Mol	Chain	Residues	Atoms					AltConf	Trace
30	BF	213	Total	C	N	O	S	0	0
			1561	1010	281	269	1		

- Molecule 31 is a protein called 60S ribosomal protein L8(A).

Mol	Chain	Residues	Atoms					AltConf	Trace
31	BG	113	Total	C	N	O	S	0	0
			844	540	144	158	2		

- Molecule 32 is a protein called 60S ribosomal protein L9(A).

Mol	Chain	Residues	Atoms					AltConf	Trace
32	BH	179	Total	C	N	O	S	0	0
			1418	896	260	259	3		

- Molecule 33 is a protein called 60S ribosomal protein L10.

Mol	Chain	Residues	Atoms					AltConf	Trace
33	BI	165	Total	C	N	O	S	0	0
			1326	834	257	228	7		

- Molecule 34 is a protein called 60S ribosomal protein L11(B).

Mol	Chain	Residues	Atoms					AltConf	Trace
34	BJ	151	Total	C	N	O	S	0	0
			1195	744	229	218	4		

- Molecule 35 is a protein called 60S ribosomal protein L12.

Mol	Chain	Residues	Atoms					AltConf	Trace
35	BK	138	Total	C	N	O	S	0	0
			1038	651	190	195	2		

- Molecule 36 is a protein called 60S ribosomal protein L15(A).

Mol	Chain	Residues	Atoms					AltConf	Trace
36	BL	192	Total	C	N	O	S	0	0
			1618	1011	340	266	1		

- Molecule 37 is a protein called 60S ribosomal protein L16(A).

Mol	Chain	Residues	Atoms					AltConf	Trace
37	BM	178	Total	C	N	O	S	0	0
			1317	845	254	217	1		

- Molecule 38 is a protein called 60S ribosomal protein L17(A).

Mol	Chain	Residues	Atoms					AltConf	Trace
38	BN	150	Total	C	N	O		0	0
			1189	742	230	217			

- Molecule 39 is a protein called 60S ribosomal protein L18(A).

Mol	Chain	Residues	Atoms					AltConf	Trace
39	BO	121	Total	C	N	O	S	0	0
			931	598	170	162	1		

- Molecule 40 is a protein called 60S ribosomal protein L19.

Mol	Chain	Residues	Atoms					AltConf	Trace
40	BP	176	Total	C	N	O		0	0
			1317	816	277	224			

- Molecule 41 is a protein called 60S ribosomal protein L21(A).

Mol	Chain	Residues	Atoms					AltConf	Trace
41	BQ	116	Total	C	N	O	S	0	0
			893	564	173	153	3		

- Molecule 42 is a protein called 60S ribosomal protein L23.

Mol	Chain	Residues	Atoms					AltConf	Trace
42	BR	131	Total	C	N	O	S	0	0
			977	614	183	173	7		

- Molecule 43 is a protein called 60S ribosomal protein L24(A).

Mol	Chain	Residues	Atoms				AltConf	Trace
43	BS	45	Total	C	N	O	0	0
			371	238	73	60		

- Molecule 44 is a protein called 60S ribosomal protein L25.

Mol	Chain	Residues	Atoms					AltConf	Trace
44	BT	80	Total	C	N	O	S	0	0
			642	411	108	121	2		

- Molecule 45 is a protein called 60S ribosomal protein L26(A).

Mol	Chain	Residues	Atoms				AltConf	Trace
45	BU	116	Total	C	N	O	0	0
			916	576	179	161		

- Molecule 46 is a protein called 60S ribosomal protein L28.

Mol	Chain	Residues	Atoms					AltConf	Trace
46	BV	142	Total	C	N	O	S	0	0
			1123	717	218	185	3		

- Molecule 47 is a protein called 60S ribosomal protein L31(A).

Mol	Chain	Residues	Atoms					AltConf	Trace
47	BW	79	Total	C	N	O	S	0	0
			663	415	135	112	1		

- Molecule 48 is a protein called 60S ribosomal protein L35.

Mol	Chain	Residues	Atoms					AltConf	Trace
48	BX	86	Total	C	N	O	S	0	0
			605	379	111	114	1		

- Molecule 49 is a protein called 60S ribosomal protein L37(A).

Mol	Chain	Residues	Atoms					AltConf	Trace
49	BY	52	Total	C	N	O	S	0	0
			403	245	85	69	4		

- Molecule 50 is a protein called 60S ribosomal protein L42.

Mol	Chain	Residues	Atoms					AltConf	Trace
50	BZ	92	Total	C	N	O	S	0	0
			749	472	151	121	5		

- Molecule 51 is a RNA chain called 5S ribosomal RNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
51	B3	113	Total	C	N	O	P	0	0
			2403	1075	429	787	112		

- Molecule 52 is a RNA chain called 5.8S ribosomal RNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
52	B4	157	Total	C	N	O	P	0	0
			3329	1490	581	1102	156		

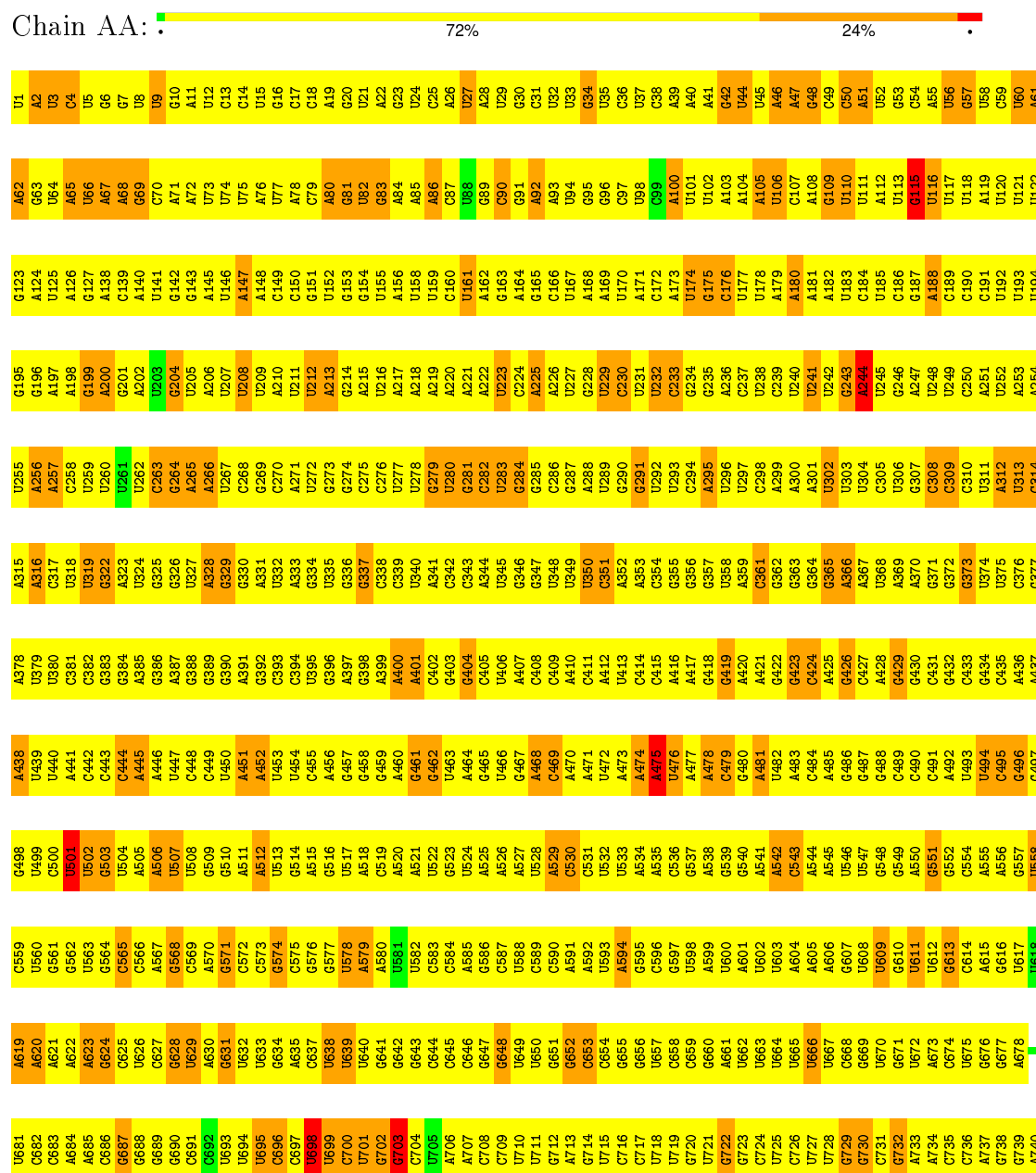
- Molecule 53 is a RNA chain called 26S ribosomal RNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
53	B5	3170	Total	C	N	O	P	0	0
			67775	30273	12178	22155	3169		

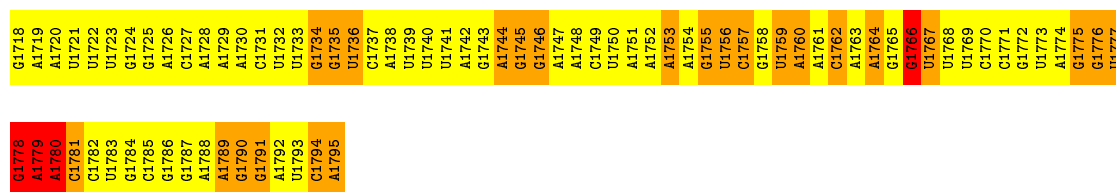
3 Residue-property plots

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. 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. 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: 18S rRNA

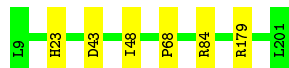


A1653	A1598	G1538	G1478	G1415	G1355	U1295	A1234	C1174	A1113	G1048	U988	A928	G868	A804	C741
U1659	G1599	G1539	C1479	G1416	C1356	G1296	U1235	G1175	U1114	U1049	C989	A929	G869	U805	U742
G1660	C1600	G1540	C1480	C1417	C1357	G1297	U1236	C1176	G1115	G1050	G990	U930	C870	A806	U743
G1661	U1601	A1481	A1481	A1418	U1358	U1298	G1237	C1177	G1116	U1051	A991	U931	G871	A807	U744
C1662	G1602	U1542	G1482	A1419	U1359	U1299	A1238	U1178	U1117	U1052	A992	A932	G872	U808	A746
G1663	G1603	A1543	C1483	U1420	U1360	U1300	G1239	A1179	C1118	U1053	G993	C933	U873	A809	C747
U1664	G1604	A1421	G1484	A1421	C1361	G1301	A1240	A1180	G1119	U1054	A994	U934	C874	G810	U748
A1665	G1605	A1422	A1485	A1422	C1362	U1302	G1241	U1181	C1120	U1055	U995	G935	C875	A811	U749
G1666	G1606	G1346	G1486	C1423	U1363	C1303	C1242	U1182	A1121	U1056	G996	C936	G876	A812	U750
U1667	U1607	C1547	U1487	A1424	G1364	U1304	U1243	U1183	A1122	G1057	A997	G937	G877	U813	G751
G1668	G1608	A1548	C1488	G1425	G1365	G1305	C1244	U1184	G1123	A1058	U998	A938	G878	A814	A752
A1669	A1609	U1549	U1489	G1426	U1366	C1306	U1245	A1185	G1124	A1059	C999	A939	G879	G815	A753
G1670	U1610	U1550	U1490	U1427	U1367	U1307	U1246	C1186	C1125	U1060	U1000	A940	C880	G816	A754
G1671	G1611	G1551	A1491	C1428	A1368	U1308	U1247	U1187	U1126	G1061	G1001	G941	A881	A817	A755
C1672	A1612	U1552	C1492	U1429	U1369	A1309	C1248	C1188	G1127	A1062	G1002	C942	U882	C818	A756
C1673	C1613	A1553	C1493	G1430	C1370	A1310	U1249	A1189	A1128	C1063	U1003	A943	C883	U820	A757
U1674	G1614	A1554	U1494	U1431	C1371	U1311	U1250	A1190	A1129	C1064	A1004	U944	G884	U821	U758
C1675	U1615	U1555	U1495	A1433	A1372	G1312	G1251	C1191	A1130	A1066	C1005	U945	G885	U822	U759
A1676	C1616	U1556	G1496	U1434	C1373	G1313	A1252	A1192	C1131	C1067	C1006	U946	U886	G823	A760
G1677	C1617	A1557	G1497	G1435	U1374	C1314	U1253	C1193	U1132	U1068	G1007	G947	U887	G824	G761
G1678	U1618	U1558	C1498	C1436	U1375	G1315	U1254	G1194	U1133	C1069	U1008	C948	U888	U825	A762
A1679	U1619	U1559	C1499	C1437	C1376	A1316	U1255	G1195	A1134	G1070	C1009	C949	U889	U826	G763
U1680	G1620	G1560	G1500	C1438	U1377	U1317	U1256	G1196	A1135	G1071	C1010	A950	C890	C827	U764
C1681	C1621	C1561	A1501	U1439	U1378	A1318	G1257	G1197	A1136	U1072	U1011	A951	A891	U828	G765
U1682	U1622	U1562	G1502	U1440	A1379	A1319	U1258	A1198	G1137	A1073	A1012	G952	A892	A829	U766
C1683	C1623	C1563	A1503	A1443	G1380	C1320	G1259	A1199	G1138	C1074	G1013	G953	U893	U830	U767
G1684	U1624	U1564	G1504	A1444	A1381	G1321	G1260	A1200	A1139	C1075	U1014	A954	U894	U831	C768
U1685	U1625	U1565	G1505	C1445	G1382	A1322	G1261	C1201	A1140	U1076	C1015	C955	G895	U832	A769
A1686	U1626	C1566	U1506	G1446	G1383	A1323	U1262	U1202	U1141	U1077	U1016	G956	U896	U833	A770
G1687	G1627	A1567	C1507	U1447	G1384	C1324	G1263	C1203	U1142	A1078	U1017	U957	C897	G834	A771
G1688	U1628	A1568	U1508	U1448	A1385	G1325	G1264	A1204	G1143	C1079	A1018	U958	G898	U835	G772
A1689	A1629	C1569	U1509	C1449	C1386	A1326	U1265	C1205	A1144	G1080	A1019	U959	U899	U836	C773
G1690	C1630	G1570	G1510	U1450	U1387	G1327	G1266	C1206	C1145	A1081	C1020	U960	G900	G837	A774
A1691	A1631	A1571	G1511	G1451	A1388	A1328	G1267	A1207	G1146	G1082	C1021	C961	U901	G838	G775
A1692	C1632	G1572	U1512	G1452	U1389	C1329	U1268	G1208	G1147	A1083	A1022	A962	U902	U839	G776
G1693	A1633	G1573	A1513	G1453	C1390	U1330	G1269	G1209	A1148	A1084	U1023	G963	G903	U840	C777
C1634	C1634	A1574	A1514	C1454	G1391	U1331	C1270	U1210	A1149	A1085	A1024	U964	A904	U841	G778
G1695	C1635	A1575	U1515	C1455	G1392	U1332	A1271	C1211	G1150	U1086	A1025	A965	A905	C842	U779
G1696	G1636	U1576	C1516	G1456	U1393	A1333	U1272	C1212	G1151	C1087	A1026	A966	A906	U843	A780
G1697	C1637	U1577	U1517	C1457	U1394	A1334	G1273	A1213	G1152	A1088	C1027	U967	U907	A844	U781
A1698	C1638	C1578	U1518	A1458	U1395	C1335	G1274	G1214	C1153	A1089	U1028	C968	U908	G845	U782
A1699	G1639	C1579	G1519	C1459	C1396	C1336	C1275	A1215	A1154	A1090	A1029	A969	C909	G846	G783
U1700	G1640	U1580	U1520	G1460	A1397	U1337	C1276	C1216	C1155	G1091	U1030	A970	U910	A847	C784
C1701	U1641	A1581	G1521	C1461	A1398	A1338	G1277	A1217	C1156	U1092	G1031	G971	U911	C848	U785
U1702	C1642	G1582	A1522	G1462	G1399	C1339	U1278	C1218	C1158	C1093	C1032	A972	G912	C849	C786
C1703	G1643	U1583	A1523	C1463	C1400	U1340	U1279	A1219	U1159	G1098	C1033	A973	G913	U850	G787
G1704	C1644	A1584	A1524	G1464	C1401	A1341	U1280	A1220	A1160	G1099	G1034	C974	A914	U851	A788
A1705	U1645	C1585	C1525	C1465	G1402	A1342	C1281	U1221	G1161	U1100	A1035	G975	U915	C852	A789
U1706	A1646	G1586	U1526	U1466	A1403	A1343	U1282	A1222	G1162	U1101	G1036	A976	U916	G853	U790
G1707	G1647	C1587	C1527	A1467	U1404	U1344	A1284	A1223	A1163	C1102	U1037	A977	U917	U854	A791
U1708	U1648	G1588	C1528	C1468	G1405	A1345	G1285	G1224	G1164	U1103	A1038	A978	A918	A855	U792
C1709	A1649	C1589	G1529	A1469	G1406	G1346	U1286	G1225	U1165	G1104	G1039	G979	U919	A856	A796
A1710	U1650	A1590	U1530	C1470	A1407	G1347	U1287	A1226	G1166	G1105	A1040	U980	U920	U857	A797
G1711	C1651	A1591	C1531	U1471	A1408	G1348	G1288	U1227	G1167	G1106	U1041	U981	G921	G858	G797
A1712	G1652	G1592	G1532	G1472	G1409	G1349	G1289	U1228	A1168	G1107	C1042	A982	A922	A859	C798
G1713	A1653	U1593	U1533	A1473	U1410	U1350	U1290	G1229	G1169	G1108	G1043	G983	A923	U860	A799
C1714	U1654	C1594	G1534	C1474	U1411	G1351	G1291	A1230	C1170	G1109	G1044	G984	G924	U864	U800
G1715	U1655	A1595	C1535	G1475	U1412	C1352	G1292	C1231	C1171	A1110	U1045	G985	A925	A865	G801
G1716	G1656	U1596	U1536	G1476	G1413	U1353	A1293	A1232	U1172	G1111	U1046	G986	C926	G866	G802
A1717	A1657	C1597	G1537	A1477	A1414	A1354	G1294	G1233	G1173	U1112	G1047	A987	U927	G867	A803



- Molecule 2: 40S ribosomal protein S0(A)

Chain AB: 97%



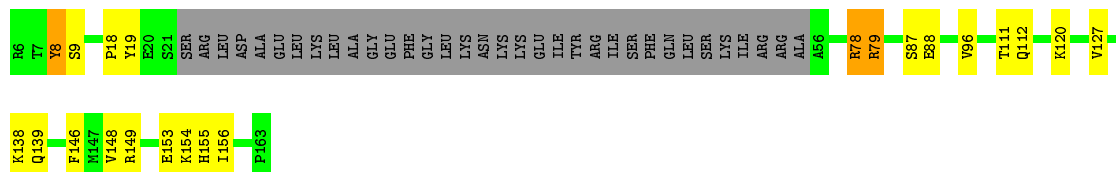
- Molecule 3: 40S ribosomal protein S3

Chain AC: 93%



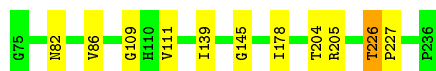
- Molecule 4: 40S ribosomal protein S9(A)

Chain AD: 65%



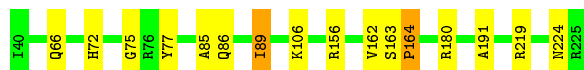
- Molecule 5: 40S ribosomal protein S2

Chain AE: 93%



- Molecule 6: 40S ribosomal protein S5

Chain AG: 91%

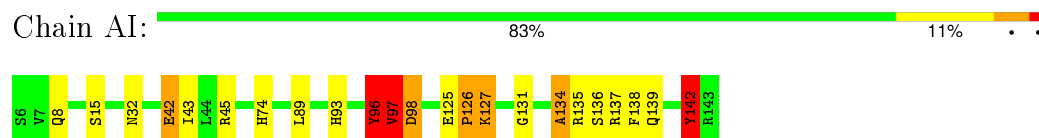


- Molecule 7: 40S ribosomal protein S22(A)

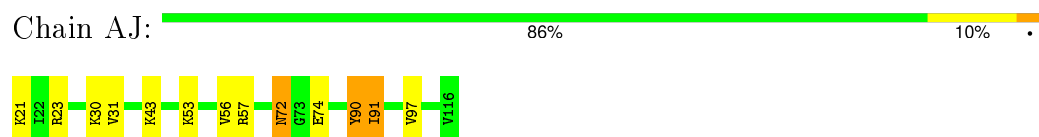
Chain AH: 89%



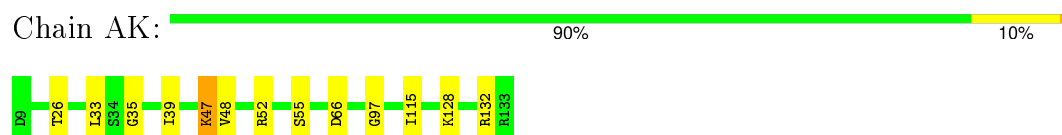
- Molecule 8: 40S ribosomal protein S16(A)



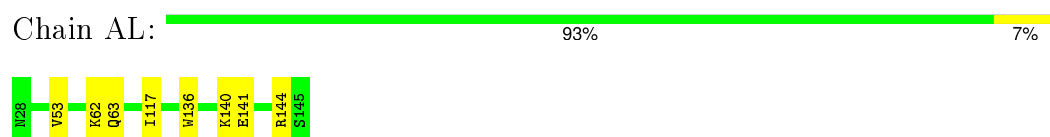
- Molecule 9: 40S ribosomal protein S20



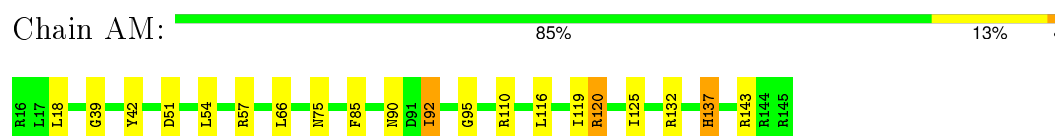
- Molecule 10: 40S ribosomal protein S14(A)



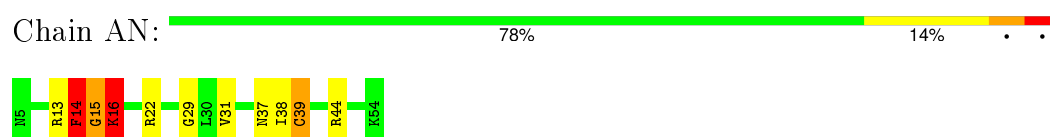
- Molecule 11: 40S ribosomal protein S23(A)



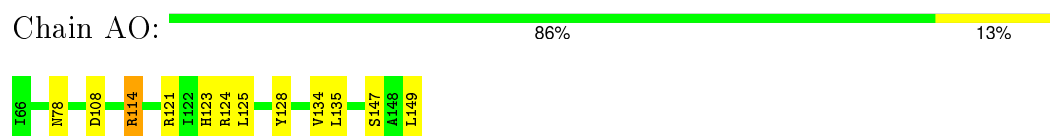
- Molecule 12: 40S ribosomal protein S18



- Molecule 13: 40S ribosomal protein S29(A)



- Molecule 14: 40S ribosomal protein S13



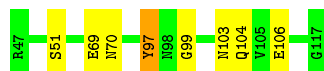
- Molecule 15: 40S ribosomal protein S11

Chain AQ:  93% 6%

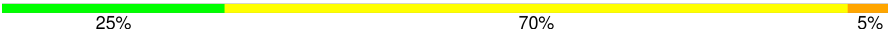


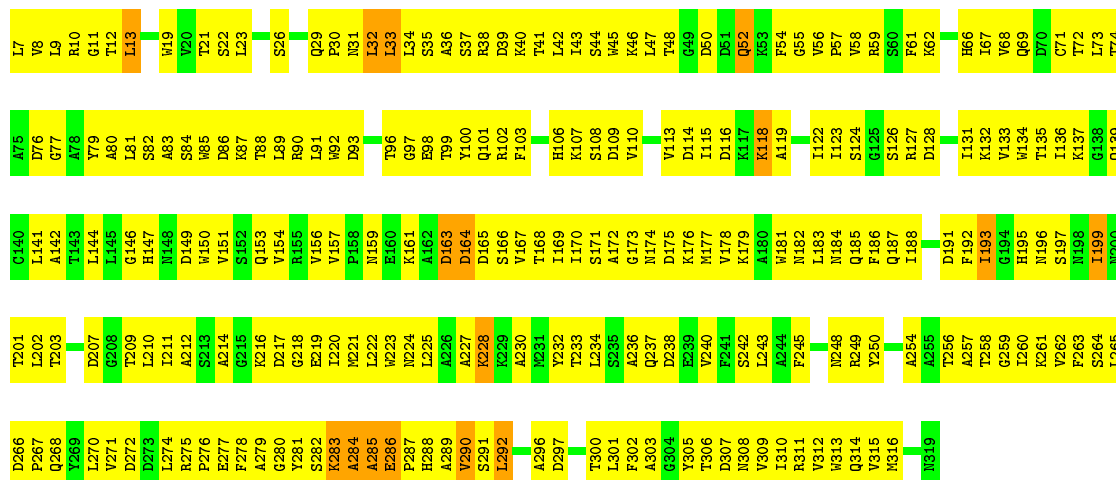
- Molecule 16: 40S ribosomal protein S15

Chain AS:  89% 10%




- Molecule 17: RACK1 protein

Chain AR:  25% 70% 5%



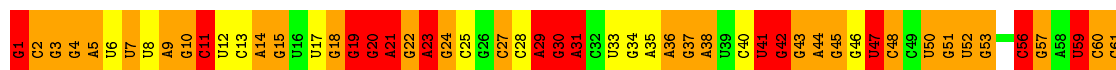
- Molecule 18: s19e protein

Chain AT:  82% 15%




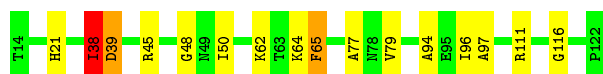
- Molecule 19: P/E tRNA

Chain A7:  11% 21% 42% 26%



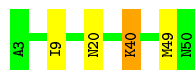
- Molecule 20: 60S ribosomal protein L32

Chain B0:  85% 12%



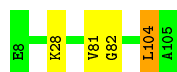
- Molecule 21: 60S ribosomal protein L39

Chain B1: 92% 6% •



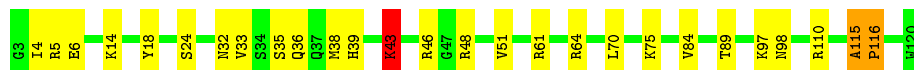
- Molecule 22: 60S ribosomal protein L30e

Chain B2: 96% ••



- Molecule 23: 60S ribosomal protein LP0

Chain B8: 77% 20% ••



- Molecule 24: 60S ribosomal protein L43

Chain B9: 82% 14% •



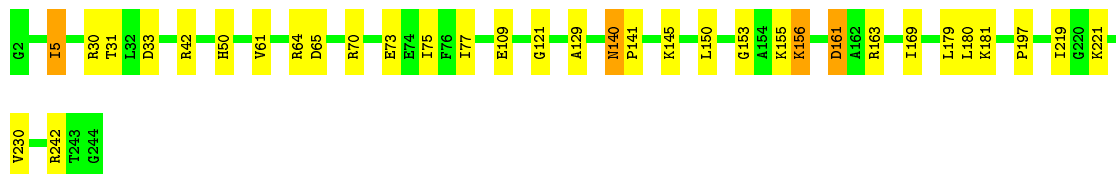
- Molecule 25: 60S ribosomal protein L1

Chain BA: 91% 8% •



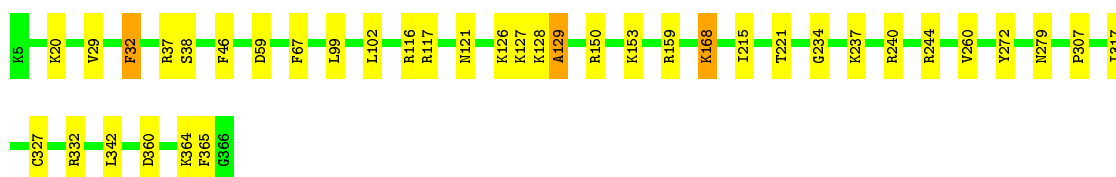
- Molecule 26: 60S ribosomal protein L2

Chain BB: 86% 12% •



- Molecule 27: 60S ribosomal protein L3

Chain BC: 90% 10% •



- Molecule 28: 60S ribosomal protein L4(B)

Chain BD: 91% 8% .



- Molecule 29: 60S ribosomal protein L5

Chain BE: 86% 13% .



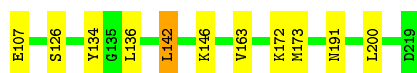
- Molecule 30: 60S ribosomal protein L7(A)

Chain BF: 93% 6% .



- Molecule 31: 60S ribosomal protein L8(A)

Chain BG: 90% 9% .



- Molecule 32: 60S ribosomal protein L9(A)

Chain BH: 88% 11% .



- Molecule 33: 60S ribosomal protein L10

Chain BI: 88% 9% .



- Molecule 34: 60S ribosomal protein L11(B)

Chain BJ: 89% 10% .



- Molecule 35: 60S ribosomal protein L12

Chain BK: 89% 10% .



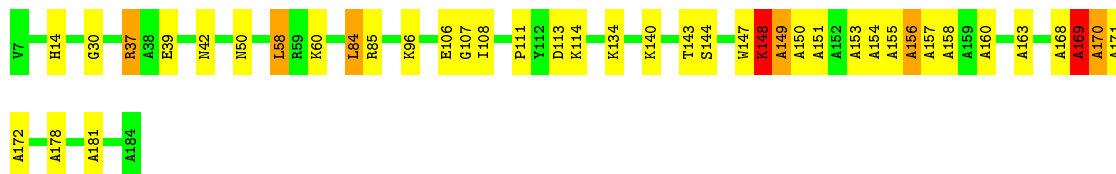
- Molecule 36: 60S ribosomal protein L15(A)

Chain BL: 90% 9% .



- Molecule 37: 60S ribosomal protein L16(A)

Chain BM: 77% 19% . .



- Molecule 38: 60S ribosomal protein L17(A)

Chain BN: 93% 7%



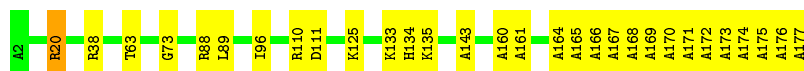
- Molecule 39: 60S ribosomal protein L18(A)

Chain BO: 94% 5% .



- Molecule 40: 60S ribosomal protein L19

Chain BP: 83% 16% .



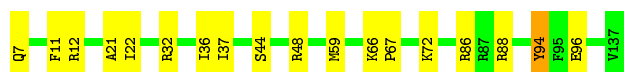
- Molecule 41: 60S ribosomal protein L21(A)

Chain BQ: 91% 9%



- Molecule 42: 60S ribosomal protein L23

Chain BR: 86% 13% .



- Molecule 43: 60S ribosomal protein L24(A)

Chain BS: 82% 16% .



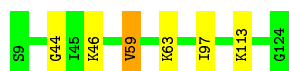
- Molecule 44: 60S ribosomal protein L25

Chain BT: 86% 13% .



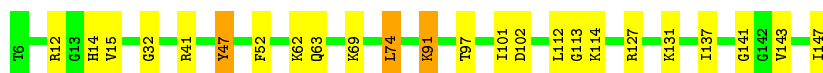
- Molecule 45: 60S ribosomal protein L26(A)

Chain BU: 95% . .



- Molecule 46: 60S ribosomal protein L28

Chain BV: 83% 15% .



- Molecule 47: 60S ribosomal protein L31(A)

Chain BW: 91% 6% .

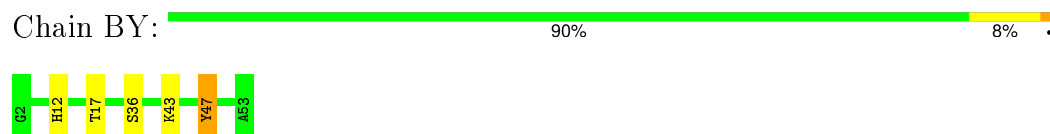


- Molecule 48: 60S ribosomal protein L35

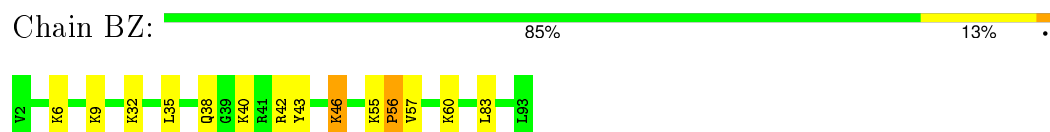
Chain BX: 86% 14%



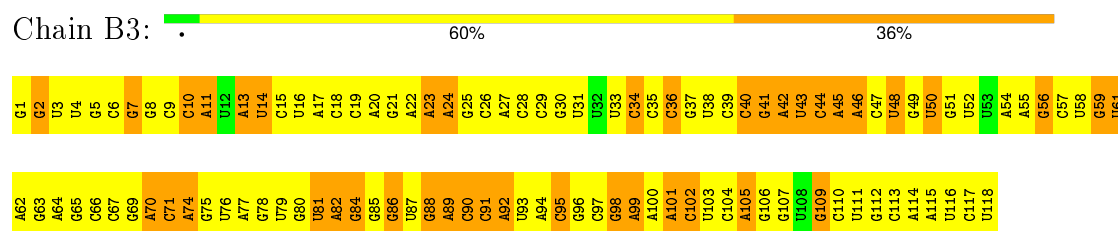
- Molecule 49: 60S ribosomal protein L37(A)



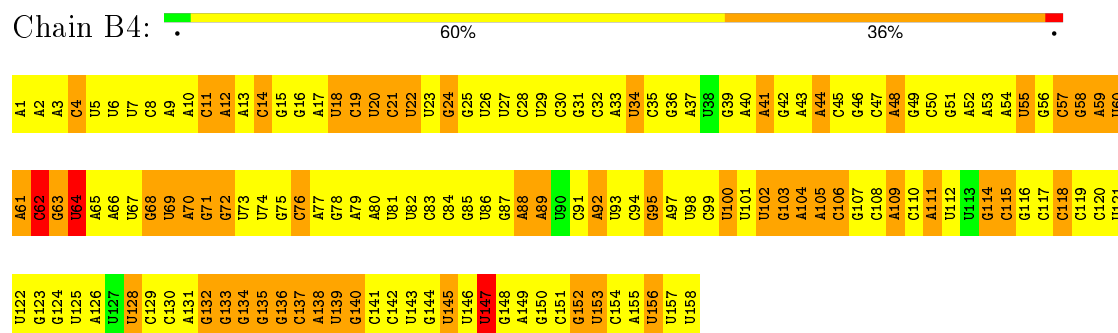
- Molecule 50: 60S ribosomal protein L42



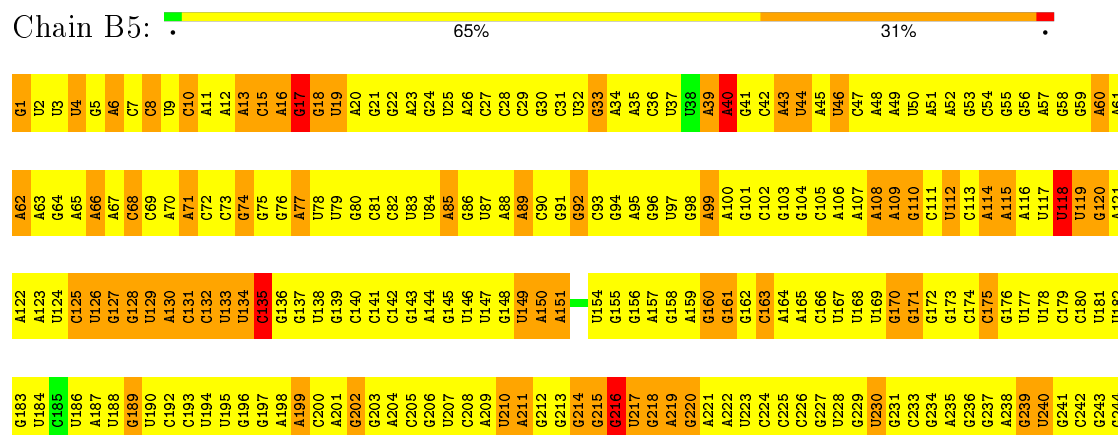
- Molecule 51: 5S ribosomal RNA



- Molecule 52: 5.8S ribosomal RNA



- Molecule 53: 26S ribosomal RNA



A1231	A1165	A1105	U1044	A974	A914	G853	C793	G733	U669	G609	U549	C489	C427	A367	A306	U245
C1232	G1166	G1106	G1045	G975	A915	G854	U794	C734	C670	G610	A550	A490	A428	A368	A307	U246
G1233	U1167	C1107	A1046	C976	G916	U855	G795	A735	U671	A611	A551	C491	U429	A369	A308	C247
G1234	U1168	U1108	A1047	U977	A917	G856	U796	A736	A672	U612	G552	U492	U430	U370	U310	U248
U1235	A1169	U1109	C1048	C978	C918	G857	U797	G737	U673	G613	U553	G493	U431	G371	U311	U249
G1236	A1170	U1110	U1049	G979	U919	A858	G798	A738	U674	G614	A554	G494	G432	A372	C312	U250
C1237	G1171	U1111	U050	U986	A920	G859	G799	G739	C675	U615	U555	G495	A433	A373	A313	C251
C1238	U1172	A1112	U1051	U987	A921	C860	G800	G740	A676	G616	U556	C496	U434	A374	U314	U252
U1239	U1173	U1113	U1052	U988	U922	C861	A801	U741	A677	G617	A557	C497	U435	A375	C315	A253
A1240	A1174	U1114	A1053	U989	C923	U862	C802	G742	U678	C618	U558	A498	A438	A376	U316	A254
C1175	C1175	U1115	A1054	A990	G924	C863	C803	G743	U679	A619	A559	A499	C439	A377	U317	A255
G1242	G1176	G1116	A1055	U991	A925	G864	C804	U744	A680	U620	G560	C500	A440	A378	A318	G256
G1243	G1177	U1117	U1056	G992	A926	U865	G805	G745	U681	A621	C561	C501	U441	C379	A319	U257
A1244	G1178	C1118	A1057	A993	C927	A866	A806	A746	U682	A622	C562	U502	G442	U380	G320	G258
A1245	A1179	U1119	U1058	G994	C928	G867	A807	A747	U683	U623	U563	C503	G443	U381	C321	C259
G1246	A1180	A1120	G1059	G995	A929	C868	U808	U748	G684	G624	U564	C504	U444	U382	U322	C260
U1247	U1181	U1121	U1060	U996	U930	G869	G809	G749	G685	G625	U565	G505	G445	G383	G323	U261
U1248	A1182	U1122	A1061	A997	G931	G870	A810	G750	G688	U626	G566	U506	U446	A384	A324	U262
G1249	C1183	U1123	A1062	A998	U932	U871	U811	A751	U689	U627	G567	U507	U447	A385	A325	C263
G1250	A1184	U1124	G1063	G999	A933	U872	G812	C752	A690	A628	U568	U508	U448	A386	U326	G264
A1251	C1185	U1125	A1064	C1000	G934	U874	G813	C753	A691	U629	U569	U509	U449	A387	A327	A265
A1252	G1186	G1126	A1065	A1002	U935	G875	U814	G754	A692	A630	U570	C510	G450	A388	U328	A266
U1253	C1187	G1127	A1066	A1003	A936	A876	G815	A755	C695	U631	A571	G511	U451	A389	U329	G267
G1254	U1188	U1128	U1067	U1004	G937	C877	A816	U756	C696	G632	A572	U512	G452	G390	G330	A268
C1255	C1189	A1129	C1068	G1005	G938	G878	A817	C757	A697	C633	C573	G513	C453	A391	G331	G269
G1256	A1190	A1130	C1069	A1006	U939	U879	C818	C758	U698	C634	U574	G514	C454	G392	C332	U270
C1257	U1191	G1131	U1070	A1007	G940	G880	U819	U759	A699	G635	C575	C515	C455	U393	G333	C271
U1258	C1192	U1132	U1071	U1008	G941	C881	A820	G760	C700	G636	C576	A516	U456	A394	C340	G272
A1259	A1193	A1133	A1072	A1009	U942	A882	G821	A761	U707	C637	C577	G517	U457	A395	G341	A273
A1260	A1200	G1134	U1073	G1010	U943	A883	G822	U762	C702	G638	A578	G518	U458	A396	A342	G274
G1261	C1201	A1135	U1074	A1011	C944	A884	C823	G763	G703	G639	G579	A519	U459	A397	G343	U275
A1262	A1202	A1136	A1075	G1012	C945	U885	C824	U764	U704	U640	C580	U520	C460	A398	A338	U276
A1263	C1203	C1137	C1076	G1013	U946	C886	U825	C765	A705	C641	U581	A521	U461	A399	C339	G277
G1264	A1204	U1138	U1077	G1014	G947	G887	G826	U766	U706	U642	G582	A522	C462	A400	C340	U278
U1265	A1205	G1139	U1078	C1017	C948	A888	A827	U767	G707	U643	G583	A523	C463	U401	G341	U279
G1266	G1206	G1140	A1079	G1018	C949	U889	U828	C768	A709	A644	A584	C524	U464	A402	A342	U280
U1267	C1207	C1141	U1080	G1019	G950	C890	U829	G769	A709	A645	A585	C525	U465	A403	U343	G281
U1268	U1208	G1142	U1081	G1020	A951	G891	A830	G770	A710	A646	C586	C526	U466	G404	A344	G282
U1269	G1209	A1143	U1082	G1021	A952	U892	G831	A771	A711	A647	U587	A527	U467	U405	G345	G283
A1270	U1210	U1144	G1083	U1022	G953	C893	G832	U772	G712	C648	G588	U528	U468	G406	C346	A284
A1271	U1211	G1145	A1084	C1023	U954	G894	U834	G773	U713	C649	A589	A529	U469	A407	G347	A285
C1272	A1212	C1146	A1085	G1024	U955	A895	U834	G774	G714	C650	G590	G530	G470	A408	A348	U286
G1273	G1213	G1147	C1086	A1025	U956	A896	G835	A775	A715	G651	U591	G531	U471	A409	A349	G287
A1274	U1214	G1148	G1087	A1026	C957	U897	A836	U776	G716	G652	G592	A532	A472	U410	C350	C288
C1275	U1215	G1149	U1088	A1027	C958	U898	A837	U777	C717	A653	A593	A533	G473	U411	A351	A289
U1276	C1216	A1150	G1089	G1029	C959	U899	G838	U778	G718	C654	A594	U534	G474	G412	A352	G290
C1277	A1217	G1090	U1090	A1030	U960	C899	C839	G779	U719	C655	A595	G535	G475	U413	G353	C291
A1278	U1218	U1151	A1091	C1031	C961	G901	C840	A780	A720	U656	U596	U536	G476	U414	U354	U292
G1279	G1219	A1153	C1092	C1032	A962	G902	A841	G781	G721	A657	G597	A537	A477	G415	A355	C293
C1280	U1220	C1154	A1093	U1033	G963	U903	G842	U782	G722	G658	A598	G538	A478	A416	C356	U294
G1281	A1221	C1155	U1094	U1034	G964	A904	A843	A783	U723	G659	C599	G539	U479	A417	A357	A295
C1282	G1222	U1095	U1095	G1035	A965	U905	G844	A784	U724	A660	G600	U540	U480	A418	G358	A296
G1283	A1223	G1157	U1096	A1036	U966	A906	G845	G785	G725	U661	U601	U541	U481	G419	U359	G297
C1284	C1224	G1097	C1037	C1037	A967	G907	A846	A786	G726	U662	A602	G542	C482	G420	G360	U298
G1285	A1225	A1159	A1098	C1038	G968	G908	A847	G787	G727	G663	A603	C543	G483	G421	A361	G299
A1286	G1226	U1099	U1039	U1039	C969	G909	A848	G788	G728	U664	G604	C544	C484	A422	U362	G300
A1287	C1227	G1161	U1100	A1040	A970	G910	C849	A789	C729	A665	U605	U545	A485	A423	G363	G301
U1288	C1228	U1162	G1101	U1041	G971	C911	U850	U790	C730	A666	C606	C546	U486	G424	G364	U302
G1289	U1229	A1163	A1102	U1042	C972	G912	U851	A791	C671	C667	A607	G547	U487	G425	A365	G303
A1290	G1230	G1164	G1104	C1043	A973	A913	U852	G792	C732	G668	A608	G548	U488	G426	A366	G304

G2377	A2317	C2196	U2135	A1921	G1860	A1800	U1737	G1674	U1607	A1546	G1483	C1423	U1361	A1291
C2378	U2318	C2197	C2136	A1922	G1861	U1801	C1738	G1675	C1608	G1547	G1485	C1424	G1362	C1292
U2379	U2319	A2198	U2137	G1923	A1862	C1802	U1739	A1676	C1609	C1548	G1486	U1425	G1363	U1293
U2380	A2320	G2199	A2138	U1924	G1863	C1803	U1740	G1677	G1610	U1549	G1487	C1426	C1364	A1294
G2381	C2321	G2261	U2200	U1925	A1864	A1804	A1741	G1678	G1611	G1488	G1489	U1427	G1365	C1296
G2382	A2322	G2201	U2140	C1926	A1865	C1805	U1742	A1679	A1612	C1551	A1428	A1366	C1296	C1296
G2383	G2323	C2202	U2141	G1927	C1866	A1806	G1743	G1680	A1613	G1552	G1429	G1423	C1297	C1297
A2384	G2324	U2203	A2142	G1928	A1867	G1807	G1744	U1681	A1614	U1553	A1491	U1430	U1368	C1298
G2385	C2325	C2204	A2143	G1929	G1868	G1808	C1745	U1682	G1615	U1554	G1492	G1431	U1369	C1299
A2386	U2326	G2206	A2144	A1930	C1869	A1809	U1746	A1683	G1616	U1555	G1493	C1432	G1370	G1300
A2387	U2327	A2207	A2145	U1931	C1870	A1810	G1747	U1684	A1617	C1556	G1494	A1433	G1371	A1301
U2388	U2328	A2208	C2146	A1932	G1871	G1811	A1748	C1685	G1618	A1557	G1495	G1434	C1372	A1302
U2389	U2329	A2147	A2147	A1933	U1872	G1812	U1749	U1686	U1622	A1558	C1497	A1435	C1373	A1303
A2390	C2330	U2209	U2148	G1934	U1873	A1813	A1750	U1687	G1623	A1559	A1498	U1436	G1374	A1304
G2391	C2331	U2210	A2149	G1935	A1874	A1814	G1751	U1688	G1624	G1560	C1499	C1437	G1375	U1305
G2392	A2332	C2212	G2150	A1936	U1875	U1815	A1752	U1689	U1626	G1561	G1500	U1438	C1376	G1306
G2393	C2333	A2213	C2151	U1937	U1876	A1816	G1753	C1690	U1627	C1562	U1501	U1439	G1377	G1307
G2394	U2334	A2214	A2152	U1938	U1877	G1817	G1754	U1691	C1628	C1563	C1502	G1440	U1378	A1308
G2395	G2335	G2215	U2153	G1939	G1878	U1818	C1755	U1692	U1629	U1564	G1441	G1441	G1379	U1309
G2396	U2336	G2216	A2154	G1940	A1879	U1819	C1756	C1693	G1630	G1565	A1504	U1442	G1380	G1310
A2397	C2337	U2217	G2155	C1941	U1880	U1820	A1757	U1694	C1631	C1505	C1505	G1443	G1381	G1311
A2398	C2338	G2218	C2156	U1942	A1881	U1821	G1758	U1695	A1632	U1567	A1506	G1444	G1382	C1312
A2399	C2339	A2219	G2157	C1943	G1882	C1822	C1759	A1696	C1633	U1568	G1507	U1445	G1383	G1313
G2400	U2340	A2220	A2158	U1944	A1883	A1823	A1760	A1697	G1634	C1508	A1508	A1446	U1384	C1314
A2401	A2341	G2221	U2159	A1945	A1884	U1824	C1761	C1698	G1635	U1570	C1509	G1447	C1385	C1316
A2402	U2342	A2222	G2160	A1946	U1885	G1825	C1762	A1699	U1636	A1571	C1510	U1448	G1387	A1317
G2403	C2343	A2223	G2161	G1947	A1886	C1826	U1763	G1700	A1637	U1572	U1511	A1449	U1388	A1318
A2404	U2344	A2224	U2162	U2102	A1887	C1827	U1764	C1701	A1638	G1573	U1512	G1450	G1389	G1319
G2405	C2345	U2225	A2163	U2103	U1888	A1828	U1765	U1702	C1639	C1574	G1513	A1451	A1390	C1320
G2406	C2346	U2226	C2164	A2104	G1889	G1829	G1766	U1703	G1640	A1575	G1514	A1452	C1391	G1321
G2407	U2347	C2227	A2166	G2105	U1890	G1830	C1767	A1704	G1641	G1576	A1515	A1453	G1392	U1322
A2408	C2348	A2228	A2167	A2106	G1892	U1831	U1768	U1705	A1642	G1577	A1516	A1454	A1393	G1323
U2409	U2349	A2229	A2168	A2107	A1893	U1832	U1769	C1706	A1643	C1578	G1517	A1455	A1394	U1324
U2410	C2350	G2230	G2169	C2108	A1894	G1833	G1770	A1707	C1644	C1579	U1518	A1456	G1395	U1325
U2411	U2351	C2231	U2170	U2109	A1895	U1834	C1771	G1709	U1646	A1580	G1519	A1457	C1396	A1326
G2412	A2352	A2232	G2171	G2110	A1896	A1835	U1772	C1710	G1646	C1581	G1520	U1458	C1397	C1327
G2413	G2353	A2233	A2172	G2111	G1897	C1836	C1773	C1711	A1647	C1582	G1521	C1459	U1398	C1328
G2414	C2354	G2234	U2173	U2112	G1898	U1837	C1774	G1712	A1648	A1583	G1521	A1460	U1329	U1329
G2415	G2355	C2235	G2174	A2113	G1899	G1838	G1775	A1714	U1649	U1584	A1524	A1461	A1401	U1330
U2416	A2356	G2236	U2175	C2114	A1900	A1839	G1776	A1715	G1650	C1585	G1525	A1462	C1402	A1331
U2417	U2357	C2237	U2176	G2115	A1901	U1840	U1777	U1716	U1651	G1586	U1526	U1463	C1403	A1332
G2418	C2358	G2238	G2177	G2116	G1902	A1841	G1778	U1717	G1652	A1587	C1527	G1464	C1333	C1333
A2419	C2359	G2239	A2178	A2117	U1903	A1842	C1779	G1718	G1653	A1588	G1528	U1405	U1334	U1334
U2420	C2360	C2179	C2179	C2118	G1904	C1843	G1780	G1719	A1654	G1590	A1529	A1466	C1335	C1335
U2421	A2361	U2241	G2180	A2119	G1905	C1844	C1781	U1720	G1655	G1591	U1530	A1467	U1336	U1336
C2422	C2362	A2242	C2181	A2120	G1906	G1845	U1782	U1721	A1656	G1592	C1531	A1468	A1337	A1337
U2423	A2363	A2243	A2182	G2121	C1907	C1846	U1783	U1722	C1657	A1593	C1532	C1469	G1409	C1338
A2424	C2364	A2244	A2183	G2122	A1908	A1847	G1784	A1723	G1658	A1594	U1533	U1470	U1410	C1339
G2425	G2365	C2245	U2184	G2123	A1909	G1848	A1787	U1724	U1659	U1595	A1534	U1471	G1340	G1340
U2426	U2366	G2246	G2185	G2124	A1910	C1849	C1788	C1725	G1660	C1596	A1535	U1472	U1341	U1341
U2427	A2367	G2247	U2186	A2125	A1911	A1850	G1789	C1726	G1661	C1597	G1536	G1473	C1342	C1342
U2428	C2368	C2248	G2187	A2126	U1912	G1851	G1790	G1727	C1665	G1598	A1537	A1474	A1343	A1343
G2429	G2369	G2249	U2188	U2127	A1913	G1852	C1791	G1728	G1666	G1599	U1538	A1475	U1415	G1344
A2430	C2370	U2189	U2189	C2128	G1914	U1853	G1793	A1729	A1667	U1600	A1539	G1476	G1345	G1345
G2431	G2371	G2251	U2190	U2129	A1915	G1854	G1794	A1731	G1668	U1601	U1540	A1477	G1346	G1346
A2432	C2372	A2252	U2191	G2130	U1916	U1855	U1795	U1732	C1669	A1602	G1541	C1478	A1418	U1347
U2433	A2373	C2253	C2192	A2131	C1917	C1856	G1796	G1733	C1670	A1603	G1542	U1479	A1419	G1357
U2434	C2374	U2254	U2193	C2132	G1918	C1857	A1797	G1734	C1671	G1604	G1543	G1480	C1420	C1358
G2435	G2375	A2255	U2194	U2133	A1858	G1735	A1798	G1735	U1672	A1481	G1421	A1481	G1422	C1359
U2436	G2376	C2256	C2195	G2134	U1920	A1859	A1799	G1736	G1673	U1606	A1545	A1482	G1422	C1360

A3347	G3285	C3225	C3164	U3104	C3043	C2983	U2923	G2863	A2803	A2740	A2679	G2618	U2558	U2497	A2437
G3346	G3286	A3226	A3165	U3105	G3044	C2984	U2924	A2864	A2804	C2741	A2680	G2619	A2559	U2498	A2438
G3349	U3287	A3227	C3166	A3106	G3045	C2985	G2925	U2865	U2806	G2742	U2681	G2620	U2560	U2499	A2439
C3350	G3288	A3228	A3167	U3107	A3046	C2986	A2926	U2866	U2807	A2743	U2682	G2621	A2561	A2500	G2440
U3351	G3289	G3229	A3168	G3108	U3047	A2987	C2927	C2867	U2808	U2744	U2683	C2622	G2562	U2501	A2441
U3352	G3290	A3169	U3169	G3109	C3048	C2988	C2928	U2868	A2808	G2745	C2684	G2563	G2503	G2503	G2442
G3353	G3291	U3231	A3170	C3110	A3049	U2989	C2929	U2869	C2809	A2746	C2685	G2624	G2504	U2504	A2443
U3354	A3292	G3232	U3171	G3111	U3050	U2990	A2930	C2870	C2810	A2747	A2686	G2625	U2505	U2505	C2444
G3355	A3294	C3233	G3172	G3112	U3051	A2991	C2931	C2871	A2811	U2748	C2687	C2627	U2506	U2506	A2445
G3356	A3295	A3234	A3173	A3113	G3052	U2992	U2932	U2872	C2812	G2749	U2688	A2628	C2507	U2507	U2446
U3357	A3296	G3235	A3174	A3114	G3053	C2993	A2933	U2873	A2813	U2750	C2689	A2629	U2508	A2447	
U3358	U3297	U3236	C3175	G3115	U3054	A2994	A2934	G2874	G2814	G2751	G2690	C2630	U2509	G2448	
A3359	G3298	U3237	C3176	G3116	U3055	A2995	U2935	U2875	G2815	U2752	U2691	U2631	U2510	A2449	
C3360	A3299	G3238	U3177	C3117	U3056	U2996	A2936	C2876	G2816	G2753	A2692	G2632	U2511	A2511	G2450
G3361	U3300	G3239	U3178	C3118	U3057	U2997	G2937	G2877	A2817	G2754	C2693	U2633	C2572	G2512	G2451
A3362	U3301	C3240	U3179	U3119	U3058	U2998	G2938	G2878	U2818	C2755	A2694	U2634	G2573	U2513	G2452
U3363	G3302	G3241	U3180	C3120	G3059	U2999	G2939	C2879	A2819	C2756	A2695	A2635	G2574	U2514	U2453
G3364	U3303	G3242	G3181	U3121	C3060	A3000	A2940	U2880	A2820	U2757	A2696	A2636	G2575	A2515	U2454
U3365	U3304	A3243	U3182	A3122	G3061	C3001	A2941	C2881	C2821	A2758	A2697	A2637	G2576	U2516	U2455
G3366	A3305	A3244	A3183	A3123	G3062	C3002	C2942	U2882	U2822	U2759	G2698	C2638	C2577	U2517	A2456
C3367	U3306	A3245	U3184	G3124	C3063	G3003	G2943	U2883	G2823	U2760	G2699	G2639	C2578	G2518	G2457
U3368	A3307	G3246	A3185	U3125	U3064	C3004	U2944	C2884	G2824	A2761	G2700	A2640	G2579	A2519	A2458
G3369	C3308	G3247	A3186	C3126	G3065	A3005	G2945	C2885	C2825	U2762	U2701	U2641	A2580	A2520	A2459
U3370	G3309	U3187	U3187	A3127	U3066	A3006	A2946	U2886	U2826	U2763	A2702	U2642	U2581	U2521	U2460
G3371	A3310	C3248	G3188	G3128	C3067	U3007	G2947	A2887	U2827	C2764	A2703	A2643	C2582	G2522	G2461
A3372	C3311	U3250	G3189	A3129	U3068	A3008	C2948	U2888	G2828	C2765	A2704	C2644	C2583	A2523	A2462
U3373	U3312	U3251	C3190	A3130	G3069	C3009	U2949	C2889	U2829	U2766	A2705	G2645	G2584	A2524	G2463
U3374	U3313	G3252	G3191	U3131	U3070	U3010	U2950	A2890	G2830	U2767	G2706	C2646	G2585	G2525	U2464
A3375	A3314	G3253	U3192	C3132	U3071	C3011	G2951	U2891	G2831	U2768	C2707	A2647	G2586	G2526	G2465
G3376	G3315	G3254	C3193	C3133	C3072	A3012	G2952	A2892	C2832	A2769	C2708	G2648	U2587	G2527	G2466
C3379	A3316	U3255	C3194	C3134	A3073	U3013	U2953	C2893	A2833	G2770	C2709	A2649	U2588	G2528	G2467
U3380	U3317	G3256	U3195	U3135	G3074	U3014	U2954	C2894	G2834	C2773	C2710	U2650	G2589	A2529	G2468
U3381	G3318	C3257	U3196	G3136	G3075	G3015	U2955	C2895	U2835	C2774	C2711	G2651	A2590	G2530	G2469
U3319	U3319	U3259	U3198	C3137	C3076	A3016	A2956	A2896	C2836	U2775	U2712	U2652	A2591	C2531	U2470
U3384	A3320	C3260	G3199	A3139	A3077	C3018	U2958	G2898	A2838	G2778	G2714	C2654	A2593	G2533	U2472
G3385	C3321	G3261	G3200	G3140	U3079	U3019	C2959	C2899	G2839	A2779	A2715	U2655	C2594	G2534	C2473
U3387	A3322	U3262	C3201	A3141	G3080	U3020	C2960	C2890	C2840	A2780	U2716	A2656	A2595	G2474	G2474
C3388	U3323	U3263	G3202	A3142	C3081	A3021	G2961	C2891	G2841	U2781	U2717	A2657	U2596	A2536	C2475
U3389	G3325	G3264	G3144	C3143	C3082	G3022	U2962	A2902	U2842	U2782	U2718	G2658	U2597	U2537	C2476
G3390	U3326	G3265	G3145	C3145	G3083	A3023	C2963	A2903	U2843	U2783	U2719	G2659	G2598	U2538	G2477
U3391	G3327	C3266	A3206	C3146	C3084	A3024	G2964	U2904	C2844	G2784	G2720	G2660	U2599	C2478	C2478
U3392	U3328	U3267	C3207	G3147	G3085	C3025	U2965	U2905	A2845	A2785	A2721	G2661	C2600	A2540	C2479
U3393	U3329	G3268	C3208	G3148	A3086	G3026	G2966	C2906	U2846	G2786	U2722	G2662	A2601	U2541	A2480
A3394	U3330	A3269	U3209	U3149	A3087	A3027	A2967	C2907	A2847	G2787	U2723	G2663	G2602	U2542	G2480
G3395	U3331	G3270	U3210	A3150	C3088	G3028	G2968	C2908	G2848	U2788	U2724	C2664	G2603	U2543	U2482
U3396	C3332	A3271	U3211	U3151	C3089	A3029	A2969	C2909	C2849	U2789	U2725	U2665	U2604	U2544	G2483
U3397	U3333	U3272	U3212	U3152	U3090	G3030	C2970	U2910	G2850	A2790	C2726	C2666	G2605	C2545	A2484
U3398	A3334	U3273	G3213	U3153	A3091	G3031	A2971	A2911	A2851	G2791	G2728	A2667	G2606	C2546	A2485
U3399	U3335	G3274	U3214	U3154	C3092	A3032	G2972	C2912	C2852	A2792	U2729	U2668	G2607	A2547	A2486
G3337	G3337	A3275	A3215	C3154	A3094	A3033	G2973	C2913	A2853	G2793	G2730	G2669	G2608	G2548	U2487
U3399	U3338	U3276	U3216	U3155	U3095	C3034	U2974	C2914	U2854	G2794	U2731	G2670	A2609	A2549	A2488
A3399	A3339	C3277	A3217	U3156	C3096	A3035	U2975	U2915	U2855	U2795	G2732	A2671	G2610	U2550	C2489
U3399	G3340	U3278	A3218	U3157	C3097	G3036	A2976	U2916	G2856	G2796	A2733	G2672	U2611	C2551	C2490
U3399	U3341	A3279	G3219	G3158	G3098	U3037	G2977	C2917	C2857	C2797	A2734	A2673	U2612	C2552	A2491
A3399	A3342	C3280	G3220	C3159	C3099	U3038	U2978	G2918	U2858	C2798	U2735	A2674	U2613	U2553	C2492
G3343	G3343	U3281	C3221	U3160	U3100	C3039	U2979	A2919	U2859	A2799	A2736	C2675	G2614	U2554	U2493
A3344	U3282	U3222	U3222	C3161	G3101	U3040	U2980	U2920	U2860	G2800	C2737	A2676	G2615	G2555	G2494
G3345	A3345	C3223	G3102	C3162	G3102	U3041	U2981	U2921	U2861	A2801	A2738	G2677	C2616	U2556	G2495
U3346	G3284	U3284	A3103	A3163	A3103	U3042	A2982	G2922	U2862	A2802	A2739	A2678	U2617	G2557	G2496

4 Experimental information

Property	Value	Source
Reconstruction method	Not provided	Depositor
Imposed symmetry	POINT, Not provided	Depositor
Number of images	Not provided	Depositor
Resolution determination method	Not provided	Depositor
CTF correction method	Not provided	Depositor
Microscope	FEI TECNAI F30	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	25	Depositor
Minimum defocus (nm)	Not provided	Depositor
Maximum defocus (nm)	Not provided	Depositor
Magnification	39000	Depositor
Image detector	KODAK SO163 FILM	Depositor

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: OMC, 5MU, OMG, H2U, YYG, 2MG, 5MC, 1MA, M2G, 7MG, PSU

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 2$	RMSZ	$\# Z > 2$
1	AA	1.91	76/41893 (0.2%)	2.21	3096/65278 (4.7%)
10	AK	0.93	0/935	1.07	0/1257
11	AL	0.88	0/920	0.98	0/1226
12	AM	0.94	0/1094	1.08	2/1468 (0.1%)
13	AN	3.66	2/427 (0.5%)	1.42	4/567 (0.7%)
14	AO	0.92	0/707	1.04	0/950
15	AQ	0.96	0/656	1.04	0/885
16	AS	0.90	0/559	1.02	2/748 (0.3%)
17	AR	0.91	1/2463 (0.0%)	1.24	7/3350 (0.2%)
18	AT	0.96	2/1118 (0.2%)	1.10	6/1498 (0.4%)
19	A7	3.12	162/1483 (10.9%)	3.79	362/2311 (15.7%)
2	AB	0.87	0/1535	0.91	0/2097
20	B0	0.91	0/898	1.03	0/1201
21	B1	1.04	0/431	1.09	0/570
22	B2	0.85	0/760	0.89	0/1020
23	B8	0.94	0/965	1.09	4/1298 (0.3%)
24	B9	0.94	0/546	1.14	3/729 (0.4%)
25	BA	0.84	0/1709	0.97	0/2295
26	BB	0.95	0/1882	1.06	1/2528 (0.0%)
27	BC	0.91	0/2953	1.05	3/3974 (0.1%)
28	BD	0.89	0/1987	0.98	4/2690 (0.1%)
29	BE	0.94	0/1956	1.11	8/2646 (0.3%)
3	AC	0.94	0/1488	1.00	0/1996
30	BF	0.85	0/1593	0.98	2/2160 (0.1%)
31	BG	0.80	0/853	1.01	0/1153
32	BH	0.89	0/1437	1.06	0/1935
33	BI	0.96	0/1352	1.12	3/1815 (0.2%)
34	BJ	0.97	0/1212	1.03	0/1622
35	BK	1.98	6/1049 (0.6%)	1.04	1/1408 (0.1%)
36	BL	1.01	0/1652	1.03	1/2211 (0.0%)
37	BM	0.97	2/1341 (0.1%)	1.33	6/1808 (0.3%)
38	BN	0.93	0/1212	1.03	4/1631 (0.2%)

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >2	RMSZ	# Z >2
39	BO	0.92	0/943	0.97	0/1274
4	AD	0.93	0/1035	1.13	5/1390 (0.4%)
40	BP	0.94	0/1334	0.98	0/1791
41	BQ	0.91	0/909	0.98	1/1216 (0.1%)
42	BR	0.92	0/992	1.13	2/1333 (0.2%)
43	BS	1.03	0/380	1.12	2/504 (0.4%)
44	BT	0.85	0/649	1.00	2/873 (0.2%)
45	BU	0.93	0/927	0.96	0/1237
46	BV	0.88	0/1152	1.07	2/1542 (0.1%)
47	BW	0.99	0/673	1.06	0/894
48	BX	0.82	0/607	1.03	2/818 (0.2%)
49	BY	0.95	0/413	1.10	2/548 (0.4%)
5	AE	0.86	0/1227	0.99	0/1663
50	BZ	0.91	0/761	1.12	3/1006 (0.3%)
51	B3	1.34	1/2686 (0.0%)	2.18	192/4184 (4.6%)
52	B4	1.35	0/3719	2.19	267/5791 (4.6%)
53	B5	1.35	3/75857 (0.0%)	2.17	5569/118271 (4.7%)
6	AG	0.91	0/1472	0.98	1/1982 (0.1%)
7	AH	0.95	2/1008 (0.2%)	1.17	2/1351 (0.1%)
8	AI	2.93	6/1106 (0.5%)	1.15	5/1481 (0.3%)
9	AJ	0.85	0/781	1.04	2/1053 (0.2%)
All	All	1.46	263/179697 (0.1%)	1.97	9578/268527 (3.6%)

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
1	AA	2	28
10	AK	0	4
12	AM	0	3
13	AN	0	2
14	AO	0	4
19	A7	0	44
2	AB	0	1
20	B0	0	1
23	B8	0	3
24	B9	0	3
26	BB	1	4
27	BC	0	1
28	BD	0	3

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Mol	Chain	#Chirality outliers	#Planarity outliers
29	BE	1	7
30	BF	1	0
31	BG	2	2
32	BH	0	3
33	BI	0	2
34	BJ	0	3
35	BK	0	2
36	BL	1	7
37	BM	0	5
38	BN	0	2
39	BO	0	1
4	AD	0	3
40	BP	0	4
42	BR	0	3
43	BS	0	1
44	BT	0	1
45	BU	0	1
46	BV	0	3
47	BW	0	2
48	BX	2	1
49	BY	0	1
5	AE	0	1
50	BZ	0	5
51	B3	1	0
52	B4	0	6
53	B5	33	112
7	AH	1	0
8	AI	0	3
9	AJ	0	3
All	All	45	285

All (263) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	AA	856	A	N3-C4	94.83	1.91	1.34
1	AA	856	A	C5-C4	83.45	1.97	1.38
1	AA	1200	A	N3-C4	80.10	1.82	1.34
1	AA	501	U	C2-N3	76.02	1.91	1.37
1	AA	856	A	C6-N1	74.78	1.87	1.35
13	AN	16	LYS	N-CA	71.90	2.90	1.46
1	AA	1200	A	C5-C4	68.16	1.86	1.38
1	AA	501	U	C4-C5	68.10	2.04	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	AA	1200	A	C6-N1	66.53	1.82	1.35
1	AA	856	A	C5-C6	65.06	1.99	1.41
1	AA	501	U	N1-C2	62.64	1.95	1.38
1	AA	501	U	N1-C6	57.45	1.89	1.38
1	AA	501	U	N3-C4	56.90	1.89	1.38
1	AA	856	A	C2-N3	56.28	1.84	1.33
1	AA	856	A	N1-C2	53.59	1.82	1.34
1	AA	1200	A	C5-C6	51.88	1.87	1.41
1	AA	1200	A	C2-N3	46.90	1.75	1.33
1	AA	1200	A	N1-C2	44.72	1.74	1.34
8	AI	96	TYR	CE1-CZ	40.35	1.91	1.38
8	AI	96	TYR	CG-CD1	39.95	1.91	1.39
8	AI	96	TYR	CE2-CZ	39.90	1.90	1.38
8	AI	96	TYR	CG-CD2	39.33	1.90	1.39
1	AA	501	U	C5-C6	36.66	1.67	1.34
8	AI	96	TYR	CD1-CE1	33.09	1.89	1.39
8	AI	96	TYR	CD2-CE2	32.79	1.88	1.39
35	BK	121	PHE	CG-CD2	28.86	1.82	1.38
35	BK	121	PHE	CG-CD1	28.18	1.81	1.38
35	BK	121	PHE	CE2-CZ	22.09	1.79	1.37
35	BK	121	PHE	CE1-CZ	21.43	1.78	1.37
35	BK	121	PHE	CD1-CE1	20.51	1.80	1.39
35	BK	121	PHE	CD2-CE2	18.75	1.76	1.39
1	AA	858	G	N3-C4	16.40	1.47	1.35
1	AA	1521	G	N3-C4	15.47	1.46	1.35
19	A7	45	G	N7-C5	14.47	1.48	1.39
1	AA	1520	U	C4'-C3'	-12.96	1.38	1.53
1	AA	864	U	C3'-C2'	-12.93	1.38	1.52
19	A7	62	A	C5-C4	-11.95	1.30	1.38
19	A7	5	A	N7-C5	11.34	1.46	1.39
19	A7	36	A	C6-N1	-11.27	1.27	1.35
1	AA	856	A	N9-C8	10.86	1.46	1.37
1	AA	864	U	C4'-C3'	10.53	1.64	1.53
19	A7	72	C	N3-C4	-10.30	1.26	1.33
19	A7	74	C	P-O5'	10.25	1.70	1.59
19	A7	57	G	N7-C5	10.20	1.45	1.39
37	BM	148	LYS	C-N	9.88	1.56	1.34
37	BM	169	ALA	C-N	9.87	1.56	1.34
19	A7	72	C	C5'-C4'	9.74	1.63	1.51
19	A7	9	A	C6-N1	-9.64	1.28	1.35
19	A7	36	A	C5-C4	9.61	1.45	1.38
19	A7	61	C	N3-C4	-9.45	1.27	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
19	A7	13	C	C4-N4	-9.32	1.25	1.33
19	A7	1	G	C6-N1	-9.26	1.33	1.39
19	A7	42	G	C2-N2	-9.14	1.25	1.34
19	A7	76	A	C5'-C4'	9.09	1.62	1.51
1	AA	1521	G	N9-C4	8.99	1.45	1.38
1	AA	1635	C	C5'-C4'	8.82	1.61	1.51
19	A7	15	G	N9-C4	8.73	1.45	1.38
19	A7	35	A	N9-C8	8.70	1.44	1.37
1	AA	858	G	C2-N2	8.64	1.43	1.34
19	A7	76	A	C5-C4	-8.59	1.32	1.38
1	AA	1521	G	C2-N2	8.54	1.43	1.34
19	A7	43	G	C2-N2	-8.53	1.26	1.34
19	A7	14	A	C6-N6	8.50	1.40	1.33
19	A7	34	OMG	O3'-P	-8.40	1.51	1.61
19	A7	57	G	N1-C2	-8.38	1.31	1.37
1	AA	864	U	O3'-P	-8.33	1.51	1.61
19	A7	41	U	P-O5'	8.30	1.68	1.59
19	A7	52	U	P-O5'	8.25	1.68	1.59
19	A7	53	G	C2'-O2'	-8.15	1.31	1.41
19	A7	9	A	C5-C4	-8.04	1.33	1.38
1	AA	1200	A	N9-C8	7.80	1.44	1.37
1	AA	858	G	N9-C4	7.75	1.44	1.38
1	AA	1516	C	P-O5'	-7.72	1.52	1.59
19	A7	13	C	C2-O2	-7.71	1.17	1.24
19	A7	25	C	C5-C6	7.66	1.40	1.34
19	A7	11	C	N3-C4	-7.63	1.28	1.33
19	A7	31	A	N7-C5	-7.61	1.34	1.39
19	A7	72	C	C4-C5	7.61	1.49	1.43
19	A7	19	G	C2-N2	-7.58	1.26	1.34
13	AN	16	LYS	CA-CB	7.56	1.70	1.53
1	AA	856	A	C8-N7	7.50	1.36	1.31
19	A7	15	G	C2-N2	-7.42	1.27	1.34
1	AA	857	U	C5'-C4'	7.37	1.60	1.51
1	AA	1584	A	C5'-C4'	7.34	1.60	1.51
19	A7	2	C	N1-C6	7.33	1.41	1.37
19	A7	14	A	N3-C4	7.33	1.39	1.34
19	A7	41	U	C2'-O2'	7.30	1.51	1.41
19	A7	60	C	C5-C6	7.28	1.40	1.34
19	A7	71	G	C5'-C4'	7.22	1.60	1.51
19	A7	64	A	C5-C6	7.17	1.47	1.41
19	A7	7	U	N1-C2	7.16	1.45	1.38
1	AA	1519	G	C4'-C3'	7.16	1.61	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
19	A7	68	U	N1-C6	7.12	1.44	1.38
19	A7	36	A	C5-C6	7.12	1.47	1.41
19	A7	22	G	N3-C4	-7.11	1.30	1.35
19	A7	71	G	C2-N2	-7.09	1.27	1.34
1	AA	1547	C	O3'-P	-7.08	1.52	1.61
19	A7	20	G	N7-C5	7.04	1.43	1.39
19	A7	23	A	C6-N1	-7.03	1.30	1.35
1	AA	856	A	C6-N6	7.00	1.39	1.33
19	A7	12	U	C4'-O4'	-7.00	1.36	1.45
1	AA	1200	A	C6-N6	7.00	1.39	1.33
19	A7	44	A	C5-C6	7.00	1.47	1.41
19	A7	44	A	N9-C4	6.98	1.42	1.37
19	A7	47	U	C5-C6	6.96	1.40	1.34
19	A7	1	G	N7-C5	6.95	1.43	1.39
19	A7	76	A	N9-C4	-6.92	1.33	1.37
19	A7	13	C	N3-C4	-6.91	1.29	1.33
19	A7	20	G	C2-N3	6.89	1.38	1.32
1	AA	859	A	P-O5'	6.88	1.66	1.59
19	A7	76	A	N7-C5	-6.84	1.35	1.39
1	AA	1583	U	C5'-C4'	6.82	1.59	1.51
19	A7	30	G	N3-C4	6.78	1.40	1.35
19	A7	45	G	C6-N1	-6.76	1.34	1.39
1	AA	1548	A	P-O5'	-6.76	1.52	1.59
19	A7	30	G	C6-N1	-6.75	1.34	1.39
1	AA	1520	U	O3'-P	-6.65	1.53	1.61
19	A7	45	G	N3-C4	6.65	1.40	1.35
1	AA	1200	A	C8-N7	6.64	1.36	1.31
19	A7	59	U	O3'-P	-6.64	1.53	1.61
19	A7	51	G	N1-C2	-6.57	1.32	1.37
19	A7	18	G	C2-N2	-6.54	1.28	1.34
1	AA	867	G	C3'-C2'	-6.54	1.45	1.52
19	A7	69	U	O3'-P	-6.52	1.53	1.61
19	A7	59	U	C2-N3	6.51	1.42	1.37
1	AA	1520	U	C3'-O3'	-6.50	1.33	1.42
1	AA	858	G	C2-N3	6.44	1.38	1.32
19	A7	23	A	C3'-C2'	-6.43	1.45	1.52
1	AA	1521	G	C5'-C4'	6.42	1.59	1.51
19	A7	35	A	C8-N7	-6.40	1.27	1.31
19	A7	36	A	N7-C5	6.39	1.43	1.39
19	A7	64	A	C2-N3	-6.37	1.27	1.33
19	A7	24	G	P-O5'	6.36	1.66	1.59
1	AA	1547	C	C3'-O3'	-6.34	1.33	1.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
19	A7	53	G	C3'-C2'	6.33	1.59	1.52
1	AA	1547	C	C4'-C3'	-6.32	1.46	1.53
18	AT	78	LYS	CA-C	6.30	1.69	1.52
19	A7	50	U	C4-O4	6.30	1.28	1.23
19	A7	4	G	C4'-O4'	-6.28	1.37	1.45
19	A7	4	G	N3-C4	6.25	1.39	1.35
19	A7	6	U	N3-C4	-6.25	1.32	1.38
19	A7	41	U	C2-O2	6.22	1.27	1.22
19	A7	56	C	N3-C4	-6.19	1.29	1.33
1	AA	1518	U	C3'-C2'	-6.18	1.46	1.52
19	A7	3	G	N3-C4	6.17	1.39	1.35
19	A7	5	A	P-O5'	-6.16	1.53	1.59
1	AA	1636	G	C3'-C2'	6.15	1.59	1.52
1	AA	1519	G	C3'-O3'	-6.15	1.33	1.42
19	A7	21	A	P-O5'	-6.12	1.53	1.59
19	A7	57	G	N9-C4	-6.12	1.33	1.38
19	A7	74	C	C4-C5	-6.12	1.38	1.43
19	A7	35	A	N7-C5	6.10	1.43	1.39
19	A7	46	7MG	O3'-P	-6.10	1.53	1.61
19	A7	50	U	P-O5'	-6.08	1.53	1.59
19	A7	41	U	C2-N3	6.05	1.42	1.37
19	A7	53	G	N7-C5	6.05	1.42	1.39
19	A7	42	G	C5'-C4'	6.02	1.58	1.51
19	A7	24	G	C2-N2	-6.00	1.28	1.34
19	A7	12	U	C5'-C4'	6.00	1.58	1.51
19	A7	20	G	C2'-C1'	6.00	1.59	1.53
19	A7	2	C	C4-N4	-5.98	1.28	1.33
1	AA	1552	U	C4'-C3'	-5.98	1.46	1.52
1	AA	1521	G	C2-N3	5.97	1.37	1.32
19	A7	61	C	C4-N4	-5.97	1.28	1.33
19	A7	66	A	C6-N1	-5.96	1.31	1.35
19	A7	21	A	O3'-P	-5.95	1.54	1.61
1	AA	1521	G	N1-C2	5.94	1.42	1.37
19	A7	31	A	C5-C4	-5.94	1.34	1.38
19	A7	76	A	C8-N7	-5.93	1.27	1.31
19	A7	51	G	C6-N1	-5.93	1.35	1.39
1	AA	1521	G	C8-N7	-5.92	1.27	1.30
19	A7	27	C	C4-N4	-5.92	1.28	1.33
51	B3	46	A	C5'-C4'	5.91	1.58	1.51
19	A7	51	G	N7-C5	-5.90	1.35	1.39
19	A7	14	A	C6-N1	-5.89	1.31	1.35
7	AH	61	ILE	CA-CB	5.89	1.68	1.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
19	A7	60	C	C2'-C1'	5.89	1.59	1.53
19	A7	69	U	O4'-C1'	5.89	1.49	1.41
19	A7	30	G	C4'-C3'	5.88	1.59	1.53
19	A7	63	C	P-O5'	-5.88	1.53	1.59
19	A7	64	A	N3-C4	5.88	1.38	1.34
17	AR	164	ASP	N-CA	-5.87	1.34	1.46
19	A7	3	G	C2'-O2'	5.87	1.49	1.41
19	A7	29	A	N3-C4	5.87	1.38	1.34
19	A7	47	U	N1-C2	5.86	1.43	1.38
19	A7	43	G	P-O5'	5.84	1.65	1.59
19	A7	67	A	C6-N6	-5.82	1.29	1.33
19	A7	19	G	C5'-C4'	5.77	1.58	1.51
1	AA	1545	A	C4'-C3'	5.76	1.59	1.53
19	A7	65	G	C8-N7	-5.75	1.27	1.30
1	AA	864	U	P-O5'	5.74	1.65	1.59
19	A7	70	C	C4-N4	-5.73	1.28	1.33
19	A7	68	U	C2-N3	5.66	1.41	1.37
7	AH	32	LYS	CA-C	5.65	1.67	1.52
19	A7	71	G	C2-N3	5.64	1.37	1.32
19	A7	69	U	C2-O2	5.63	1.27	1.22
1	AA	797	G	P-O5'	-5.63	1.54	1.59
53	B5	2229	A	C5'-C4'	5.61	1.58	1.51
19	A7	47	U	C4-O4	-5.61	1.19	1.23
1	AA	1519	G	C3'-C2'	-5.60	1.46	1.52
19	A7	67	A	C2'-O2'	5.57	1.48	1.41
19	A7	47	U	C1'-N1	5.55	1.57	1.48
19	A7	57	G	C5-C6	5.55	1.48	1.42
19	A7	42	G	N7-C5	5.52	1.42	1.39
19	A7	48	C	C3'-C2'	5.51	1.59	1.52
19	A7	57	G	C8-N7	5.50	1.34	1.30
1	AA	1635	C	C4'-C3'	5.49	1.59	1.53
1	AA	858	G	N1-C2	5.49	1.42	1.37
19	A7	5	A	C5-C4	-5.49	1.34	1.38
1	AA	859	A	C5'-C4'	5.48	1.57	1.51
19	A7	69	U	C4-O4	-5.46	1.19	1.23
19	A7	11	C	C2-O2	-5.43	1.19	1.24
1	AA	858	G	C6-N1	5.41	1.43	1.39
19	A7	56	C	N1-C6	-5.40	1.33	1.37
19	A7	73	A	C8-N7	5.40	1.35	1.31
19	A7	28	C	C2'-O2'	-5.39	1.34	1.41
19	A7	44	A	C4'-O4'	-5.39	1.38	1.45
1	AA	502	U	C5'-C4'	5.39	1.57	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	AA	1546	G	P-O5'	5.39	1.65	1.59
19	A7	60	C	O3'-P	5.37	1.67	1.61
19	A7	3	G	C5'-C4'	5.32	1.57	1.51
1	AA	1612	A	C5'-C4'	5.29	1.57	1.51
19	A7	64	A	C4'-C3'	5.29	1.58	1.53
19	A7	72	C	C3'-C2'	-5.28	1.47	1.52
1	AA	1515	U	O3'-P	-5.26	1.54	1.61
19	A7	3	G	N7-C5	-5.25	1.36	1.39
19	A7	3	G	C4'-O4'	5.25	1.52	1.45
19	A7	63	C	C5-C6	-5.25	1.30	1.34
19	A7	53	G	N1-C2	-5.22	1.33	1.37
19	A7	52	U	C4'-O4'	-5.22	1.38	1.45
1	AA	1637	C	O3'-P	-5.21	1.54	1.61
19	A7	41	U	C5-C6	5.21	1.38	1.34
19	A7	11	C	N1-C2	5.21	1.45	1.40
19	A7	12	U	O4'-C1'	5.19	1.48	1.41
1	AA	1520	U	C5'-C4'	-5.18	1.45	1.51
53	B5	1421	G	C5'-C4'	5.18	1.57	1.51
18	AT	79	LEU	N-CA	5.18	1.56	1.46
19	A7	59	U	C5-C6	5.17	1.38	1.34
19	A7	66	A	C5'-C4'	5.17	1.57	1.51
19	A7	48	C	C5-C6	-5.16	1.30	1.34
19	A7	56	C	C5-C6	5.16	1.38	1.34
1	AA	1521	G	O4'-C1'	5.16	1.48	1.41
19	A7	19	G	C5-C6	5.16	1.47	1.42
19	A7	33	U	C2'-O2'	5.16	1.48	1.41
19	A7	51	G	C3'-C2'	5.15	1.58	1.52
19	A7	48	C	O4'-C1'	5.14	1.48	1.41
19	A7	1	G	O3'-P	-5.13	1.54	1.61
19	A7	63	C	O3'-P	-5.12	1.55	1.61
19	A7	62	A	C2'-C1'	-5.12	1.47	1.53
19	A7	6	U	C4'-O4'	-5.12	1.38	1.45
1	AA	1518	U	C3'-O3'	-5.11	1.35	1.42
19	A7	35	A	N3-C4	5.10	1.38	1.34
19	A7	71	G	C5-C6	5.10	1.47	1.42
1	AA	796	A	O3'-P	-5.09	1.55	1.61
19	A7	37	YYG	O3'-P	-5.08	1.55	1.61
1	AA	1520	U	P-O5'	-5.08	1.54	1.59
19	A7	6	U	C5'-C4'	5.08	1.57	1.51
19	A7	30	G	N7-C5	5.07	1.42	1.39
19	A7	53	G	N3-C4	-5.06	1.31	1.35
19	A7	28	C	C3'-O3'	5.03	1.49	1.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
19	A7	15	G	C5-C6	5.02	1.47	1.42
1	AA	799	A	C5'-C4'	5.02	1.57	1.51
53	B5	2623	G	C2-N3	5.00	1.36	1.32

All (9578) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1545	A	O5'-P-OP1	-67.13	30.14	110.70
1	AA	864	U	P-O3'-C3'	-44.04	66.85	119.70
1	AA	1516	C	O5'-P-OP2	-30.31	74.33	110.70
1	AA	1213	A	O5'-P-OP1	-23.91	82.01	110.70
1	AA	1519	G	O3'-P-O5'	23.73	149.08	104.00
1	AA	1213	A	O5'-P-OP2	-23.37	82.65	110.70
1	AA	1520	U	O3'-P-O5'	23.19	148.06	104.00
1	AA	1521	G	N9-C4-C5	-22.98	96.21	105.40
1	AA	1516	C	O5'-P-OP1	22.56	137.77	110.70
1	AA	856	A	N7-C8-N9	21.87	124.73	113.80
1	AA	1579	C	P-O3'-C3'	21.67	145.70	119.70
37	BM	148	LYS	O-C-N	-20.85	89.34	122.70
1	AA	1547	C	P-O3'-C3'	-20.45	95.16	119.70
37	BM	169	ALA	O-C-N	-19.62	91.30	122.70
1	AA	1521	G	C4-C5-N7	19.32	118.53	110.80
1	AA	1548	A	O5'-P-OP1	19.25	133.80	110.70
1	AA	866	G	O5'-P-OP2	-18.88	88.05	110.70
19	A7	75	C	N3-C4-C5	18.68	129.37	121.90
1	AA	856	A	P-O3'-C3'	18.65	142.08	119.70
19	A7	36	A	N7-C8-N9	18.59	123.09	113.80
1	AA	867	G	OP1-P-OP2	-18.33	92.11	119.60
1	AA	865	A	O5'-P-OP2	-18.23	88.82	110.70
1	AA	865	A	P-O3'-C3'	-18.12	97.95	119.70
1	AA	858	G	C1'-O4'-C4'	-17.43	95.96	109.90
1	AA	1520	U	P-O3'-C3'	-17.41	98.80	119.70
1	AA	858	G	O4'-C4'-C3'	-17.05	86.95	104.00
1	AA	1535	C	P-O3'-C3'	17.02	140.12	119.70
1	AA	852	C	P-O3'-C3'	16.93	140.01	119.70
1	AA	1520	U	C4'-C3'-O3'	-16.51	74.72	109.40
1	AA	1196	G	P-O3'-C3'	16.48	139.48	119.70
53	B5	3143	C	P-O3'-C3'	16.18	139.12	119.70
19	A7	76	A	N1-C2-N3	-16.08	121.26	129.30
1	AA	1547	C	C4'-C3'-C2'	16.08	118.68	102.60
1	AA	1200	A	N7-C8-N9	16.04	121.82	113.80
1	AA	748	U	P-O3'-C3'	16.03	138.94	119.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	856	A	C4-C5-N7	-15.97	102.72	110.70
53	B5	2787	G	P-O3'-C3'	15.89	138.77	119.70
1	AA	858	G	O4'-C1'-N9	15.82	120.86	108.20
1	AA	856	A	N9-C4-C5	-15.71	99.52	105.80
1	AA	1520	U	C3'-C2'-C1'	-15.63	88.99	101.50
19	A7	36	A	C5-N7-C8	-15.53	96.14	103.90
1	AA	1318	A	P-O3'-C3'	15.52	138.33	119.70
1	AA	1635	C	P-O3'-C3'	-15.51	101.09	119.70
1	AA	652	G	P-O3'-C3'	15.30	138.06	119.70
1	AA	1543	A	P-O3'-C3'	-15.13	101.54	119.70
52	B4	152	G	P-O3'-C3'	15.00	137.70	119.70
19	A7	75	C	C2-N3-C4	-14.77	112.52	119.90
53	B5	1582	C	P-O3'-C3'	14.74	137.39	119.70
19	A7	76	A	C5-C6-N1	14.73	125.06	117.70
19	A7	9	A	C5-C6-N1	14.66	125.03	117.70
53	B5	475	G	P-O3'-C3'	14.64	137.26	119.70
53	B5	1163	A	P-O3'-C3'	14.56	137.18	119.70
53	B5	1302	A	P-O3'-C3'	14.53	137.13	119.70
52	B4	19	C	P-O3'-C3'	14.52	137.13	119.70
53	B5	638	C	P-O3'-C3'	14.50	137.10	119.70
1	AA	1521	G	P-O3'-C3'	-14.48	102.33	119.70
19	A7	31	A	C8-N9-C4	-14.44	100.02	105.80
53	B5	239	G	P-O3'-C3'	14.42	137.01	119.70
53	B5	712	G	P-O3'-C3'	14.35	136.92	119.70
19	A7	5	A	N1-C6-N6	-14.28	110.03	118.60
19	A7	76	A	C2-N3-C4	14.21	117.71	110.60
1	AA	1213	A	OP1-P-OP2	14.19	140.88	119.60
1	AA	1560	G	P-O3'-C3'	14.14	136.67	119.70
19	A7	44	A	N1-C6-N6	-14.14	110.11	118.60
53	B5	3160	U	P-O3'-C3'	14.13	136.65	119.70
37	BM	148	LYS	CA-C-N	-14.12	86.14	117.20
53	B5	1590	G	P-O3'-C3'	14.11	136.63	119.70
53	B5	2571	U	P-O3'-C3'	14.07	136.59	119.70
1	AA	1200	A	C4-C5-N7	-14.07	103.67	110.70
53	B5	1737	U	P-O3'-C3'	14.04	136.55	119.70
53	B5	1626	U	P-O3'-C3'	14.03	136.54	119.70
1	AA	1098	G	P-O3'-C3'	14.00	136.50	119.70
1	AA	1520	U	C2'-C3'-O3'	-13.96	78.80	109.50
1	AA	1488	C	P-O3'-C3'	13.95	136.44	119.70
1	AA	1639	C	P-O3'-C3'	13.94	136.42	119.70
1	AA	1055	U	P-O3'-C3'	13.93	136.42	119.70
1	AA	229	U	P-O3'-C3'	13.93	136.41	119.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
51	B3	23	A	P-O3'-C3'	13.93	136.41	119.70
1	AA	1675	C	P-O3'-C3'	13.91	136.39	119.70
19	A7	62	A	N1-C6-N6	-13.90	110.26	118.60
1	AA	1518	U	OP1-P-OP2	-13.87	98.80	119.60
1	AA	1429	U	P-O3'-C3'	13.80	136.26	119.70
1	AA	1759	U	P-O3'-C3'	13.74	136.19	119.70
53	B5	3384	U	P-O3'-C3'	13.72	136.17	119.70
19	A7	36	A	C4-C5-C6	-13.71	110.15	117.00
1	AA	1478	G	P-O3'-C3'	13.67	136.11	119.70
1	AA	1133	U	P-O3'-C3'	13.65	136.09	119.70
53	B5	3083	G	P-O3'-C3'	13.61	136.04	119.70
53	B5	3219	U	P-O3'-C3'	13.61	136.03	119.70
1	AA	759	U	P-O3'-C3'	13.59	136.00	119.70
1	AA	212	U	P-O3'-C3'	13.54	135.95	119.70
53	B5	2988	C	P-O3'-C3'	13.52	135.93	119.70
19	A7	53	G	C5-N7-C8	-13.51	97.55	104.30
53	B5	1033	U	P-O3'-C3'	13.50	135.90	119.70
19	A7	47	U	O4'-C1'-N1	13.49	118.99	108.20
53	B5	3049	A	P-O3'-C3'	13.48	135.88	119.70
1	AA	1629	A	P-O3'-C3'	-13.43	103.59	119.70
1	AA	1521	G	N3-C4-N9	13.42	134.05	126.00
18	AT	78	LYS	CA-C-O	-13.40	91.95	120.10
1	AA	1281	C	P-O3'-C3'	13.40	135.78	119.70
19	A7	76	A	C4-C5-C6	-13.39	110.31	117.00
53	B5	2913	C	P-O3'-C3'	13.38	135.76	119.70
1	AA	1544	G	O5'-P-OP2	-13.37	93.67	105.70
53	B5	801	A	P-O3'-C3'	13.37	135.74	119.70
19	A7	45	G	C6-C5-N7	13.35	138.41	130.40
1	AA	461	G	P-O3'-C3'	13.34	135.71	119.70
53	B5	2459	A	P-O3'-C3'	13.28	135.63	119.70
1	AA	66	U	P-O3'-C3'	13.26	135.61	119.70
1	AA	1636	G	C2'-C3'-O3'	13.23	138.62	109.50
1	AA	770	A	P-O3'-C3'	13.21	135.55	119.70
1	AA	922	A	P-O3'-C3'	13.21	135.55	119.70
53	B5	2538	U	P-O3'-C3'	13.20	135.54	119.70
52	B4	102	U	P-O3'-C3'	13.19	135.52	119.70
19	A7	9	A	N1-C6-N6	-13.18	110.69	118.60
53	B5	1084	A	P-O3'-C3'	13.17	135.50	119.70
1	AA	56	U	P-O3'-C3'	13.13	135.46	119.70
1	AA	199	G	P-O3'-C3'	13.12	135.44	119.70
1	AA	1216	C	P-O3'-C3'	13.11	135.43	119.70
53	B5	3138	U	P-O3'-C3'	13.11	135.43	119.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	866	G	O5'-P-OP1	-13.10	93.91	105.70
19	A7	64	A	N1-C6-N6	-13.03	110.78	118.60
19	A7	1	G	C5-C6-O6	-13.01	120.79	128.60
53	B5	215	G	P-O3'-C3'	12.97	135.26	119.70
1	AA	856	A	N3-C4-N9	12.94	137.75	127.40
1	AA	365	G	P-O3'-C3'	12.93	135.22	119.70
53	B5	3141	A	P-O3'-C3'	12.91	135.19	119.70
1	AA	175	G	P-O3'-C3'	12.89	135.17	119.70
1	AA	867	G	O5'-P-OP1	12.85	126.12	110.70
1	AA	1200	A	N9-C4-C5	-12.83	100.67	105.80
17	AR	163	ASP	C-N-CA	-12.80	89.69	121.70
19	A7	27	C	N3-C2-O2	-12.79	112.95	121.90
1	AA	1200	A	N1-C6-N6	12.77	126.26	118.60
1	AA	501	U	C2-N3-C4	-12.76	119.34	127.00
51	B3	91	C	P-O3'-C3'	12.76	135.02	119.70
53	B5	1549	U	P-O3'-C3'	12.76	135.02	119.70
1	AA	1789	A	P-O3'-C3'	12.72	134.96	119.70
1	AA	1457	C	P-O3'-C3'	12.71	134.95	119.70
1	AA	1181	U	P-O3'-C3'	12.68	134.91	119.70
37	BM	169	ALA	CA-C-N	-12.68	89.31	117.20
53	B5	2654	C	P-O3'-C3'	12.60	134.82	119.70
1	AA	438	A	P-O3'-C3'	12.56	134.77	119.70
1	AA	1551	G	OP1-P-OP2	-12.53	100.80	119.60
53	B5	1390	A	P-O3'-C3'	12.52	134.72	119.70
1	AA	1779	A	P-O3'-C3'	12.51	134.71	119.70
53	B5	3310	A	N1-C6-N6	12.48	126.09	118.60
53	B5	52	A	N1-C6-N6	12.47	126.08	118.60
1	AA	1635	C	C2'-C3'-O3'	12.44	136.87	109.50
52	B4	69	U	P-O3'-C3'	12.34	134.51	119.70
1	AA	350	U	P-O3'-C3'	12.32	134.48	119.70
1	AA	1636	G	O5'-P-OP2	-12.31	94.62	105.70
53	B5	372	A	P-O3'-C3'	12.20	134.34	119.70
19	A7	6	U	C5-C6-N1	-12.20	116.60	122.70
1	AA	1144	A	N1-C6-N6	12.19	125.91	118.60
53	B5	1643	A	P-O3'-C3'	12.18	134.31	119.70
37	BM	169	ALA	C-N-CA	-12.17	91.28	121.70
1	AA	864	U	N1-C1'-C2'	-12.12	98.25	114.00
19	A7	21	A	O4'-C1'-N9	12.11	117.89	108.20
19	A7	2	C	N3-C4-N4	-12.09	109.54	118.00
1	AA	1238	A	N1-C6-N6	12.08	125.85	118.60
19	A7	45	G	C5-C6-N1	12.08	117.54	111.50
19	A7	28	C	O4'-C1'-N1	12.07	117.85	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1212	C	O3'-P-O5'	12.05	126.89	104.00
19	A7	8	U	N3-C4-O4	11.97	127.78	119.40
53	B5	920	A	N1-C6-N6	11.97	125.78	118.60
19	A7	5	A	N1-C2-N3	-11.91	123.35	129.30
53	B5	134	U	P-O3'-C3'	11.88	133.95	119.70
13	AN	16	LYS	N-CA-CB	11.87	131.97	110.60
53	B5	2177	G	P-O3'-C3'	11.87	133.94	119.70
53	B5	3330	A	P-O3'-C3'	11.86	133.93	119.70
1	AA	1200	A	N3-C4-N9	11.86	136.88	127.40
52	B4	21	C	P-O3'-C3'	11.83	133.90	119.70
13	AN	15	GLY	C-N-CA	11.80	151.21	121.70
53	B5	2314	U	P-O3'-C3'	11.75	133.80	119.70
1	AA	1547	C	C3'-C2'-C1'	-11.75	92.10	101.50
19	A7	71	G	C2-N3-C4	11.73	117.77	111.90
19	A7	76	A	N1-C6-N6	-11.72	111.57	118.60
53	B5	2373	A	P-O3'-C3'	11.68	133.72	119.70
1	AA	1585	A	N1-C6-N6	11.66	125.60	118.60
19	A7	62	A	C5-C6-N1	11.66	123.53	117.70
1	AA	1265	U	P-O3'-C3'	11.65	133.69	119.70
1	AA	1520	U	OP2-P-O3'	-11.63	79.61	105.20
53	B5	810	A	N1-C6-N6	11.60	125.56	118.60
7	AH	32	LYS	CA-C-O	-11.59	95.77	120.10
53	B5	2474	G	P-O3'-C3'	11.59	133.60	119.70
53	B5	1915	A	N1-C6-N6	11.52	125.51	118.60
19	A7	33	U	O4'-C1'-N1	11.51	117.41	108.20
53	B5	2833	A	N1-C6-N6	11.49	125.50	118.60
19	A7	64	A	C4-C5-C6	-11.49	111.26	117.00
53	B5	2520	A	P-O3'-C3'	11.48	133.48	119.70
19	A7	59	U	C5-C4-O4	11.48	132.79	125.90
1	AA	1212	C	OP1-P-O3'	-11.47	79.96	105.20
53	B5	3086	A	N1-C6-N6	11.46	125.48	118.60
1	AA	1517	U	OP1-P-OP2	-11.46	102.41	119.60
53	B5	620	U	P-O3'-C3'	11.45	133.44	119.70
53	B5	791	A	N1-C6-N6	11.43	125.46	118.60
1	AA	1521	G	OP1-P-OP2	-11.43	102.46	119.60
53	B5	3016	A	N1-C6-N6	11.43	125.45	118.60
1	AA	1520	U	OP1-P-O3'	-11.42	80.08	105.20
53	B5	219	A	N1-C6-N6	11.38	125.43	118.60
1	AA	1521	G	O4'-C4'-C3'	-11.35	92.65	104.00
53	B5	2231	C	C2-N1-C1'	11.35	131.28	118.80
53	B5	820	A	N1-C6-N6	11.35	125.41	118.60
19	A7	30	G	C5-C6-N1	11.34	117.17	111.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
19	A7	27	C	C2-N3-C4	-11.32	114.24	119.90
53	B5	2748	A	N1-C6-N6	11.32	125.39	118.60
1	AA	856	A	C2-N3-C4	11.31	116.25	110.60
53	B5	1271	A	N1-C6-N6	11.29	125.38	118.60
1	AA	444	C	P-O3'-C3'	11.27	133.22	119.70
1	AA	1553	A	OP1-P-OP2	-11.24	102.74	119.60
53	B5	997	A	N1-C6-N6	11.23	125.34	118.60
1	AA	1521	G	N3-C2-N2	11.22	127.75	119.90
1	AA	1122	A	N1-C6-N6	11.22	125.33	118.60
1	AA	1788	A	N1-C6-N6	11.21	125.33	118.60
53	B5	2172	A	N1-C6-N6	11.21	125.33	118.60
19	A7	56	C	C2-N3-C4	11.20	125.50	119.90
53	B5	965	A	N1-C6-N6	11.20	125.32	118.60
53	B5	2386	A	N1-C6-N6	11.20	125.32	118.60
1	AA	100	A	N1-C6-N6	11.18	125.31	118.60
53	B5	2671	A	N1-C6-N6	11.18	125.31	118.60
53	B5	3049	A	N1-C6-N6	11.18	125.31	118.60
53	B5	628	A	N1-C6-N6	11.18	125.31	118.60
53	B5	277	G	P-O3'-C3'	11.17	133.11	119.70
53	B5	1813	A	N1-C6-N6	11.16	125.30	118.60
53	B5	433	A	N1-C6-N6	11.16	125.30	118.60
1	AA	685	A	N1-C6-N6	11.15	125.29	118.60
53	B5	2295	A	N1-C6-N6	11.14	125.29	118.60
53	B5	608	A	N1-C6-N6	11.14	125.28	118.60
19	A7	9	A	N9-C4-C5	11.13	110.25	105.80
1	AA	929	A	N1-C6-N6	11.13	125.28	118.60
53	B5	3322	A	N1-C6-N6	11.12	125.27	118.60
53	B5	2811	A	N1-C6-N6	11.11	125.27	118.60
53	B5	2897	A	N1-C6-N6	11.10	125.26	118.60
53	B5	847	A	N1-C6-N6	11.10	125.26	118.60
53	B5	1456	A	N1-C6-N6	11.10	125.26	118.60
52	B4	92	A	N1-C6-N6	11.09	125.25	118.60
53	B5	119	U	P-O3'-C3'	11.08	132.99	119.70
53	B5	1048	A	N1-C6-N6	11.07	125.24	118.60
53	B5	2447	A	N1-C6-N6	11.07	125.24	118.60
53	B5	1818	U	P-O3'-C3'	11.06	132.97	119.70
53	B5	3385	U	P-O3'-C3'	11.06	132.97	119.70
1	AA	1338	A	N1-C6-N6	11.05	125.23	118.60
53	B5	1491	A	N1-C6-N6	11.05	125.23	118.60
53	B5	318	A	N1-C6-N6	11.05	125.23	118.60
53	B5	926	A	N1-C6-N6	11.05	125.23	118.60
1	AA	867	G	O5'-P-OP2	11.04	123.95	110.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	1679	A	N1-C6-N6	11.03	125.22	118.60
53	B5	827	A	N1-C6-N6	11.03	125.22	118.60
53	B5	1904	C	O4'-C1'-N1	11.02	117.02	108.20
53	B5	2480	A	N1-C6-N6	11.02	125.21	118.60
53	B5	1165	A	N1-C6-N6	11.02	125.21	118.60
1	AA	1212	C	OP2-P-O3'	-11.00	81.00	105.20
1	AA	1159	U	P-O3'-C3'	11.00	132.90	119.70
1	AA	1654	U	P-O3'-C3'	10.98	132.88	119.70
1	AA	1226	A	N1-C6-N6	10.97	125.18	118.60
53	B5	1933	A	N1-C6-N6	10.96	125.18	118.60
1	AA	62	A	N1-C6-N6	10.96	125.17	118.60
53	B5	2279	A	N1-C6-N6	10.95	125.17	118.60
1	AA	1234	A	N1-C6-N6	10.94	125.17	118.60
53	B5	817	A	N1-C6-N6	10.92	125.15	118.60
1	AA	1491	A	N1-C6-N6	10.91	125.15	118.60
53	B5	13	A	N1-C6-N6	10.91	125.15	118.60
1	AA	1518	U	O5'-P-OP1	10.91	123.79	110.70
19	A7	30	G	N1-C6-O6	-10.90	113.36	119.90
53	B5	828	A	N1-C6-N6	10.90	125.14	118.60
1	AA	1473	A	N1-C6-N6	10.89	125.14	118.60
53	B5	26	A	N1-C6-N6	10.89	125.14	118.60
53	B5	2149	A	N1-C6-N6	10.89	125.14	118.60
53	B5	2796	G	P-O3'-C3'	10.89	132.77	119.70
51	B3	55	A	N1-C6-N6	10.89	125.14	118.60
53	B5	3320	A	N1-C6-N6	10.89	125.13	118.60
1	AA	756	A	N1-C6-N6	10.88	125.12	118.60
53	B5	324	A	N1-C6-N6	10.87	125.12	118.60
53	B5	2224	A	N1-C6-N6	10.87	125.12	118.60
52	B4	149	A	N1-C6-N6	10.85	125.11	118.60
53	B5	705	A	N1-C6-N6	10.85	125.11	118.60
19	A7	68	U	N3-C2-O2	-10.84	114.61	122.20
1	AA	1519	G	OP1-P-O3'	-10.83	81.38	105.20
53	B5	2926	A	N1-C6-N6	10.83	125.10	118.60
53	B5	2956	A	N1-C6-N6	10.83	125.10	118.60
19	A7	7	U	O4'-C1'-N1	10.82	116.86	108.20
53	B5	1715	A	N1-C6-N6	10.82	125.09	118.60
53	B5	2419	A	N1-C6-N6	10.82	125.09	118.60
53	B5	1529	A	N1-C6-N6	10.82	125.09	118.60
53	B5	89	A	N1-C6-N6	10.81	125.09	118.60
53	B5	2387	A	N1-C6-N6	10.80	125.08	118.60
53	B5	2515	A	N1-C6-N6	10.79	125.08	118.60
1	AA	68	A	N1-C6-N6	10.79	125.07	118.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
52	B4	44	A	N1-C6-N6	10.79	125.07	118.60
1	AA	799	A	O5'-P-OP2	-10.78	96.00	105.70
53	B5	1643	A	N1-C6-N6	10.78	125.07	118.60
53	B5	2691	A	N1-C6-N6	10.78	125.07	118.60
19	A7	5	A	C4-C5-C6	-10.77	111.62	117.00
53	B5	2215	A	N1-C6-N6	10.77	125.06	118.60
1	AA	1022	A	N1-C6-N6	10.77	125.06	118.60
1	AA	855	A	N1-C6-N6	10.76	125.06	118.60
1	AA	417	A	N1-C6-N6	10.76	125.05	118.60
1	AA	445	A	N1-C6-N6	10.75	125.05	118.60
53	B5	2902	A	N1-C6-N6	10.75	125.05	118.60
1	AA	1185	A	N1-C6-N6	10.74	125.05	118.60
1	AA	41	A	N1-C6-N6	10.73	125.04	118.60
19	A7	66	A	N1-C6-N6	-10.73	112.16	118.60
53	B5	1203	A	N1-C6-N6	10.72	125.03	118.60
1	AA	456	A	N1-C6-N6	10.72	125.03	118.60
53	B5	2847	A	N1-C6-N6	10.71	125.02	118.60
1	AA	804	A	N1-C6-N6	10.70	125.02	118.60
53	B5	2609	A	N1-C6-N6	10.70	125.02	118.60
53	B5	1080	A	N1-C6-N6	10.70	125.02	118.60
19	A7	1	G	C6-N1-C2	-10.69	118.69	125.10
53	B5	808	A	N1-C6-N6	10.69	125.01	118.60
19	A7	44	A	C5'-C4'-O4'	10.68	121.92	109.10
53	B5	1079	A	P-O3'-C3'	10.67	132.51	119.70
53	B5	3279	A	N1-C6-N6	10.66	125.00	118.60
1	AA	1730	A	N1-C6-N6	10.64	124.99	118.60
52	B4	43	A	N1-C6-N6	10.64	124.98	118.60
53	B5	836	A	N1-C6-N6	10.64	124.98	118.60
1	AA	1085	A	N1-C6-N6	10.63	124.98	118.60
53	B5	2229	A	N1-C6-N6	10.63	124.98	118.60
53	B5	2838	A	N1-C6-N6	10.63	124.98	118.60
53	B5	1064	A	N1-C6-N6	10.63	124.98	118.60
53	B5	1504	A	N1-C6-N6	10.63	124.98	118.60
53	B5	2703	A	N1-C6-N6	10.63	124.98	118.60
19	A7	69	U	N1-C2-N3	10.63	121.28	114.90
53	B5	784	A	N1-C6-N6	10.62	124.97	118.60
53	B5	1332	A	N1-C6-N6	10.62	124.97	118.60
53	B5	1750	A	N1-C6-N6	10.62	124.97	118.60
53	B5	3362	A	N1-C6-N6	10.62	124.97	118.60
53	B5	439	C	O4'-C1'-N1	10.61	116.69	108.20
53	B5	2139	A	N1-C6-N6	10.61	124.97	118.60
1	AA	188	A	N1-C6-N6	10.60	124.96	118.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	1112	A	N1-C6-N6	10.60	124.96	118.60
53	B5	254	A	N1-C6-N6	10.59	124.95	118.60
53	B5	1291	A	N1-C6-N6	10.59	124.96	118.60
53	B5	2326	A	N1-C6-N6	10.59	124.96	118.60
53	B5	3218	A	N1-C6-N6	10.59	124.96	118.60
53	B5	3245	A	N1-C6-N6	10.59	124.95	118.60
1	AA	1520	U	O4'-C1'-N1	-10.59	99.73	108.20
53	B5	974	A	N1-C6-N6	10.58	124.95	118.60
1	AA	951	A	N1-C6-N6	10.58	124.95	118.60
1	AA	1792	A	N1-C6-N6	10.58	124.95	118.60
53	B5	1027	A	N1-C6-N6	10.57	124.94	118.60
53	B5	1120	A	N1-C6-N6	10.57	124.94	118.60
53	B5	2837	A	N1-C6-N6	10.55	124.93	118.60
53	B5	3073	A	N1-C6-N6	10.55	124.93	118.60
53	B5	2143	A	N1-C6-N6	10.55	124.93	118.60
53	B5	1003	A	N1-C6-N6	10.54	124.92	118.60
53	B5	319	A	N1-C6-N6	10.54	124.92	118.60
53	B5	917	A	N1-C6-N6	10.54	124.92	118.60
53	B5	1804	A	N1-C6-N6	10.53	124.92	118.60
53	B5	1446	A	N1-C6-N6	10.52	124.92	118.60
1	AA	39	A	N1-C6-N6	10.52	124.91	118.60
1	AA	858	G	C5'-C4'-C3'	10.52	132.83	116.00
53	B5	386	A	N1-C6-N6	10.52	124.91	118.60
19	A7	53	G	O4'-C4'-C3'	10.51	114.51	104.00
53	B5	70	A	N1-C6-N6	10.51	124.91	118.60
53	B5	2958	A	N1-C6-N6	10.51	124.90	118.60
1	AA	1521	G	O4'-C1'-N9	10.50	116.60	108.20
19	A7	31	A	N9-C4-C5	10.50	110.00	105.80
1	AA	615	A	N1-C6-N6	10.50	124.90	118.60
1	AA	1700	A	N1-C6-N6	10.50	124.90	118.60
53	B5	1245	A	N1-C6-N6	10.50	124.90	118.60
53	B5	607	A	N1-C6-N6	10.49	124.90	118.60
53	B5	2491	A	N1-C6-N6	10.49	124.90	118.60
53	B5	2656	A	N1-C6-N6	10.49	124.90	118.60
53	B5	3113	A	P-O3'-C3'	10.49	132.29	119.70
53	B5	3323	A	N1-C6-N6	10.49	124.90	118.60
53	B5	1893	A	N1-C6-N6	10.48	124.89	118.60
53	B5	1787	A	N1-C6-N6	10.48	124.89	118.60
53	B5	2674	A	N1-C6-N6	10.48	124.89	118.60
53	B5	2679	A	N1-C6-N6	10.48	124.89	118.60
53	B5	1418	A	N1-C6-N6	10.48	124.89	118.60
53	B5	77	A	N1-C6-N6	10.47	124.89	118.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	710	A	N1-C6-N6	10.47	124.89	118.60
53	B5	3027	A	N1-C6-N6	10.47	124.88	118.60
53	B5	3305	A	N1-C6-N6	10.47	124.89	118.60
53	B5	100	A	N1-C6-N6	10.47	124.88	118.60
53	B5	2120	A	N1-C6-N6	10.47	124.88	118.60
1	AA	1524	A	N1-C6-N6	10.46	124.88	118.60
1	AA	962	A	N1-C6-N6	10.46	124.88	118.60
53	B5	2636	A	N1-C6-N6	10.46	124.88	118.60
53	B5	3168	A	N1-C6-N6	10.46	124.87	118.60
1	AA	1519	G	P-O3'-C3'	10.45	132.24	119.70
1	AA	171	A	N1-C6-N6	10.45	124.87	118.60
1	AA	521	A	N1-C6-N6	10.45	124.87	118.60
1	AA	219	A	N1-C6-N6	10.45	124.87	118.60
53	B5	1602	A	N1-C6-N6	10.45	124.87	118.60
1	AA	954	A	N1-C6-N6	10.44	124.87	118.60
1	AA	1760	A	N1-C6-N6	10.45	124.87	118.60
53	B5	201	A	N1-C6-N6	10.45	124.87	118.60
53	B5	2892	A	N1-C6-N6	10.44	124.87	118.60
53	B5	12	A	N1-C6-N6	10.44	124.86	118.60
53	B5	1287	A	N1-C6-N6	10.44	124.86	118.60
1	AA	939	A	N1-C6-N6	10.44	124.86	118.60
53	B5	3227	A	N1-C6-N6	10.44	124.86	118.60
1	AA	865	A	O3'-P-O5'	10.43	123.82	104.00
51	B3	92	A	N1-C6-N6	10.43	124.86	118.60
53	B5	2367	A	N1-C6-N6	10.43	124.86	118.60
1	AA	1217	A	N1-C6-N6	10.42	124.85	118.60
53	B5	2188	A	N1-C6-N6	10.41	124.85	118.60
19	A7	5	A	C8-N9-C4	10.41	109.96	105.80
53	B5	1381	A	N1-C6-N6	10.41	124.84	118.60
53	B5	2178	A	N1-C6-N6	10.40	124.84	118.60
53	B5	1603	A	N1-C6-N6	10.39	124.84	118.60
53	B5	761	A	N1-C6-N6	10.39	124.83	118.60
53	B5	2903	A	N1-C6-N6	10.39	124.83	118.60
1	AA	1129	A	N1-C6-N6	10.39	124.83	118.60
53	B5	3046	A	N1-C6-N6	10.38	124.83	118.60
1	AA	1026	A	N1-C6-N6	10.38	124.83	118.60
53	B5	1575	A	N1-C6-N6	10.38	124.83	118.60
53	B5	352	A	N1-C6-N6	10.38	124.83	118.60
1	AA	202	A	N1-C6-N6	10.37	124.82	118.60
1	AA	707	A	N1-C6-N6	10.38	124.83	118.60
53	B5	375	A	N1-C6-N6	10.38	124.83	118.60
53	B5	690	A	N1-C6-N6	10.37	124.82	118.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	775	A	N1-C6-N6	10.37	124.82	118.60
53	B5	1025	A	N1-C6-N6	10.37	124.82	118.60
1	AA	485	A	N1-C6-N6	10.37	124.82	118.60
53	B5	67	A	N1-C6-N6	10.36	124.81	118.60
53	B5	3039	C	P-O3'-C3'	10.36	132.13	119.70
1	AA	86	A	N1-C6-N6	10.36	124.81	118.60
53	B5	2851	A	N1-C6-N6	10.35	124.81	118.60
1	AA	914	A	N1-C6-N6	10.34	124.81	118.60
52	B4	40	A	N1-C6-N6	10.34	124.80	118.60
53	B5	409	A	N1-C6-N6	10.34	124.80	118.60
53	B5	3299	A	N1-C6-N6	10.34	124.80	118.60
53	B5	1079	A	N1-C6-N6	10.33	124.80	118.60
1	AA	928	A	N1-C6-N6	10.33	124.80	118.60
1	AA	1554	A	N1-C6-N6	10.33	124.80	118.60
53	B5	1676	A	N1-C6-N6	10.33	124.80	118.60
1	AA	1328	A	N1-C6-N6	10.33	124.80	118.60
1	AA	1134	A	P-O3'-C3'	10.33	132.09	119.70
53	B5	602	A	N1-C6-N6	10.33	124.80	118.60
53	B5	2304	C	O4'-C1'-N1	10.33	116.46	108.20
53	B5	377	A	N1-C6-N6	10.32	124.79	118.60
53	B5	2445	A	N1-C6-N6	10.32	124.79	118.60
53	B5	222	A	N1-C6-N6	10.32	124.79	118.60
19	A7	60	C	C6-N1-C2	-10.32	116.17	120.30
53	B5	2348	A	N1-C6-N6	10.32	124.79	118.60
51	B3	24	A	N1-C6-N6	10.32	124.79	118.60
53	B5	2113	A	N1-C6-N6	10.32	124.79	118.60
1	AA	1422	A	N1-C6-N6	10.31	124.79	118.60
53	B5	830	A	N1-C6-N6	10.31	124.79	118.60
51	B3	54	A	N1-C6-N6	10.31	124.79	118.60
53	B5	783	A	N1-C6-N6	10.31	124.78	118.60
53	B5	1231	A	N1-C6-N6	10.31	124.78	118.60
53	B5	2256	A	N1-C6-N6	10.31	124.78	118.60
53	B5	2195	C	O4'-C1'-N1	10.30	116.44	108.20
53	B5	2321	A	N1-C6-N6	10.30	124.78	118.60
53	B5	744	A	N1-C6-N6	10.30	124.78	118.60
53	B5	2357	A	N1-C6-N6	10.30	124.78	118.60
51	B3	101	A	N1-C6-N6	10.29	124.78	118.60
53	B5	2145	A	N1-C6-N6	10.29	124.78	118.60
1	AA	940	A	N1-C6-N6	10.29	124.78	118.60
53	B5	1133	A	N1-C6-N6	10.29	124.77	118.60
53	B5	621	A	P-O3'-C3'	10.29	132.05	119.70
53	B5	268	A	N1-C6-N6	10.29	124.77	118.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	3186	A	N1-C6-N6	10.28	124.77	118.60
53	B5	3347	A	N1-C6-N6	10.29	124.77	118.60
53	B5	1075	A	N1-C6-N6	10.28	124.77	118.60
53	B5	2312	A	N1-C6-N6	10.28	124.77	118.60
1	AA	1761	A	N1-C6-N6	10.28	124.77	118.60
52	B4	10	A	N1-C6-N6	10.28	124.77	118.60
53	B5	2341	A	N1-C6-N6	10.28	124.77	118.60
1	AA	734	A	N1-C6-N6	10.27	124.76	118.60
53	B5	1697	A	N1-C6-N6	10.27	124.76	118.60
1	AA	1754	A	N1-C6-N6	10.27	124.76	118.60
53	B5	3170	A	N1-C6-N6	10.27	124.76	118.60
1	AA	1547	C	C2'-C3'-O3'	-10.27	86.91	109.50
1	AA	1354	A	N1-C6-N6	10.27	124.76	118.60
1	AA	244	A	N1-C6-N6	10.26	124.76	118.60
51	B3	70	A	N1-C6-N6	10.26	124.76	118.60
52	B4	52	A	N1-C6-N6	10.26	124.76	118.60
53	B5	1559	A	N1-C6-N6	10.26	124.75	118.60
53	B5	619	A	N1-C6-N6	10.26	124.75	118.60
1	AA	856	A	C6-C5-N7	10.25	139.47	132.30
1	AA	1160	A	N1-C6-N6	10.25	124.75	118.60
1	AA	1612	A	O4'-C1'-N9	10.25	116.40	108.20
53	B5	417	A	N1-C6-N6	10.25	124.75	118.60
53	B5	1741	A	N1-C6-N6	10.24	124.75	118.60
53	B5	2456	A	N1-C6-N6	10.24	124.75	118.60
53	B5	2705	A	P-O3'-C3'	10.24	131.99	119.70
1	AA	148	A	N1-C6-N6	10.24	124.75	118.60
1	AA	470	A	N1-C6-N6	10.24	124.74	118.60
1	AA	341	A	N1-C6-N6	10.24	124.74	118.60
1	AA	1168	A	N1-C6-N6	10.24	124.74	118.60
1	AA	1541	A	N1-C6-N6	10.24	124.74	118.60
1	AA	162	A	N1-C6-N6	10.23	124.74	118.60
1	AA	1653	A	N1-C6-N6	10.23	124.74	118.60
53	B5	1163	A	N1-C6-N6	10.23	124.74	118.60
53	B5	1696	A	N1-C6-N6	10.23	124.74	118.60
53	B5	2368	A	N1-C6-N6	10.23	124.74	118.60
53	B5	354	U	P-O3'-C3'	10.23	131.98	119.70
53	B5	2523	A	N1-C6-N6	10.23	124.74	118.60
53	B5	2640	A	N1-C6-N6	10.23	124.74	118.60
1	AA	22	A	N1-C6-N6	10.23	124.74	118.60
52	B4	131	A	N1-C6-N6	10.23	124.74	118.60
53	B5	2167	A	N1-C6-N6	10.23	124.74	118.60
53	B5	71	A	N1-C6-N6	10.22	124.73	118.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	67	A	N1-C6-N6	10.22	124.73	118.60
53	B5	1566	A	N1-C6-N6	10.22	124.73	118.60
19	A7	9	A	O4'-C1'-N9	10.21	116.37	108.20
53	B5	973	A	N1-C6-N6	10.21	124.73	118.60
53	B5	1153	A	N1-C6-N6	10.21	124.73	118.60
53	B5	2946	A	N1-C6-N6	10.21	124.73	118.60
19	A7	44	A	C8-N9-C4	-10.21	101.72	105.80
1	AA	905	A	N1-C6-N6	10.21	124.72	118.60
53	B5	2130	G	N1-C6-O6	10.21	126.03	119.90
1	AA	621	A	N1-C6-N6	10.21	124.72	118.60
53	B5	1886	A	N1-C6-N6	10.21	124.72	118.60
19	A7	11	C	N3-C4-C5	10.20	125.98	121.90
53	B5	348	A	N1-C6-N6	10.20	124.72	118.60
1	AA	1636	G	P-O3'-C3'	10.20	131.94	119.70
53	B5	387	A	N1-C6-N6	10.20	124.72	118.60
53	B5	1910	A	N1-C6-N6	10.20	124.72	118.60
53	B5	2779	A	N1-C6-N6	10.20	124.72	118.60
4	AD	8	TYR	CB-CG-CD1	-10.20	114.88	121.00
53	B5	1435	A	N1-C6-N6	10.19	124.72	118.60
53	B5	2817	A	N1-C6-N6	10.20	124.72	118.60
1	AA	1485	A	N1-C6-N6	10.19	124.72	118.60
53	B5	884	A	N1-C6-N6	10.19	124.71	118.60
52	B4	155	A	N1-C6-N6	10.19	124.71	118.60
53	B5	308	A	N1-C6-N6	10.19	124.71	118.60
53	B5	2317	A	N1-C6-N6	10.19	124.71	118.60
53	B5	585	A	N1-C6-N6	10.18	124.71	118.60
53	B5	1159	A	N1-C6-N6	10.18	124.71	118.60
53	B5	2539	A	N1-C6-N6	10.18	124.71	118.60
53	B5	2511	A	N1-C6-N6	10.18	124.71	118.60
1	AA	247	A	N1-C6-N6	10.18	124.71	118.60
53	B5	408	A	N1-C6-N6	10.18	124.71	118.60
1	AA	771	A	N1-C6-N6	10.17	124.70	118.60
1	AA	1140	A	N1-C6-N6	10.17	124.70	118.60
53	B5	711	A	N1-C6-N6	10.17	124.70	118.60
1	AA	1319	A	N1-C6-N6	10.17	124.70	118.60
53	B5	925	A	N1-C6-N6	10.17	124.70	118.60
1	AA	370	A	N1-C6-N6	10.17	124.70	118.60
1	AA	1196	G	O4'-C1'-N9	10.16	116.33	108.20
1	AA	1310	A	N1-C6-N6	10.16	124.70	118.60
53	B5	157	A	N1-C6-N6	10.16	124.70	118.60
53	B5	1193	A	N1-C6-N6	10.16	124.70	118.60
53	B5	265	A	N1-C6-N6	10.16	124.70	118.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	1707	A	N1-C6-N6	10.16	124.70	118.60
53	B5	611	A	N1-C6-N6	10.16	124.69	118.60
53	B5	2569	A	N1-C6-N6	10.16	124.69	118.60
1	AA	1385	A	N1-C6-N6	10.15	124.69	118.60
53	B5	1062	A	N1-C6-N6	10.15	124.69	118.60
53	B5	1704	A	N1-C6-N6	10.15	124.69	118.60
52	B4	109	A	N1-C6-N6	10.15	124.69	118.60
53	B5	402	A	N1-C6-N6	10.15	124.69	118.60
53	B5	2443	A	N1-C6-N6	10.15	124.69	118.60
1	AA	1720	A	N1-C6-N6	10.14	124.69	118.60
1	AA	437	A	N1-C6-N6	10.14	124.68	118.60
19	A7	25	C	N3-C4-C5	10.14	125.96	121.90
53	B5	114	A	N1-C6-N6	10.14	124.68	118.60
53	B5	1683	A	N1-C6-N6	10.14	124.68	118.60
1	AA	965	A	N1-C6-N6	10.14	124.68	118.60
52	B4	53	A	N1-C6-N6	10.14	124.68	118.60
1	AA	301	A	N1-C6-N6	10.13	124.68	118.60
1	AA	1753	A	N1-C6-N6	10.13	124.68	118.60
53	B5	65	A	N1-C6-N6	10.13	124.68	118.60
52	B4	60	U	O4'-C1'-N1	10.13	116.30	108.20
1	AA	505	A	N1-C6-N6	10.12	124.67	118.60
1	AA	1591	A	N1-C6-N6	10.12	124.67	118.60
1	AA	859	A	O4'-C1'-N9	10.12	116.30	108.20
53	B5	2649	A	N1-C6-N6	10.12	124.67	118.60
1	AA	391	A	N1-C6-N6	10.12	124.67	118.60
1	AA	1692	A	N1-C6-N6	10.12	124.67	118.60
19	A7	36	A	O4'-C1'-N9	10.12	116.30	108.20
53	B5	1085	A	N1-C6-N6	10.12	124.67	118.60
53	B5	1273	A	N1-C6-N6	10.12	124.67	118.60
53	B5	3127	A	N1-C6-N6	10.12	124.67	118.60
53	B5	2716	U	O4'-C1'-N1	10.11	116.29	108.20
53	B5	1200	A	P-O3'-C3'	10.11	131.83	119.70
1	AA	1584	A	N1-C6-N6	10.11	124.67	118.60
52	B4	89	A	N1-C6-N6	10.11	124.67	118.60
53	B5	1847	A	N1-C6-N6	10.11	124.67	118.60
53	B5	1921	A	N1-C6-N6	10.11	124.66	118.60
53	B5	2540	A	N1-C6-N6	10.10	124.66	118.60
53	B5	1909	A	N1-C6-N6	10.10	124.66	118.60
53	B5	2931	C	O4'-C1'-N1	10.10	116.28	108.20
1	AA	1199	A	N1-C6-N6	10.10	124.66	118.60
53	B5	418	A	N1-C6-N6	10.10	124.66	118.60
53	B5	3174	A	N1-C6-N6	10.10	124.66	118.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	238	A	N1-C6-N6	10.09	124.66	118.60
1	AA	1744	A	N1-C6-N6	10.09	124.66	118.60
53	B5	1946	A	N1-C6-N6	10.09	124.66	118.60
53	B5	677	A	N1-C6-N6	10.09	124.65	118.60
53	B5	2255	A	N1-C6-N6	10.09	124.66	118.60
53	B5	1879	A	N1-C6-N6	10.09	124.65	118.60
53	B5	3336	A	N1-C6-N6	10.09	124.65	118.60
53	B5	1647	A	N1-C6-N6	10.08	124.65	118.60
53	B5	2390	A	N1-C6-N6	10.08	124.65	118.60
53	B5	2910	A	N1-C6-N6	10.08	124.65	118.60
53	B5	2432	A	N1-C6-N6	10.08	124.65	118.60
53	B5	1835	A	N1-C6-N6	10.08	124.65	118.60
53	B5	2559	A	N1-C6-N6	10.08	124.65	118.60
1	AA	400	A	N1-C6-N6	10.08	124.65	118.60
53	B5	3234	A	N1-C6-N6	10.08	124.64	118.60
1	AA	966	A	N1-C6-N6	10.07	124.64	118.60
53	B5	385	A	N1-C6-N6	10.07	124.64	118.60
53	B5	1509	A	N1-C6-N6	10.07	124.64	118.60
1	AA	605	A	N1-C6-N6	10.07	124.64	118.60
53	B5	1343	A	N1-C6-N6	10.07	124.64	118.60
53	B5	1605	A	N1-C6-N6	10.07	124.64	118.60
1	AA	1318	A	N1-C6-N6	10.06	124.64	118.60
53	B5	2358	A	N1-C6-N6	10.06	124.64	118.60
1	AA	1230	A	N1-C6-N6	10.06	124.63	118.60
53	B5	2919	A	N1-C6-N6	10.06	124.63	118.60
1	AA	1139	A	N1-C6-N6	10.05	124.63	118.60
1	AA	1154	A	N1-C6-N6	10.05	124.63	118.60
53	B5	952	A	N1-C6-N6	10.05	124.63	118.60
53	B5	1881	A	N1-C6-N6	10.05	124.63	118.60
53	B5	1290	A	N1-C6-N6	10.05	124.63	118.60
1	AA	399	A	N1-C6-N6	10.05	124.63	118.60
1	AA	1398	A	N1-C6-N6	10.05	124.63	118.60
53	B5	1757	A	N1-C6-N6	10.05	124.63	118.60
53	B5	2736	A	N1-C6-N6	10.05	124.63	118.60
1	AA	1081	A	N1-C6-N6	10.04	124.63	118.60
53	B5	416	A	N1-C6-N6	10.04	124.63	118.60
53	B5	1093	A	N1-C6-N6	10.05	124.63	118.60
53	B5	2590	A	N1-C6-N6	10.05	124.63	118.60
53	B5	1539	A	N1-C6-N6	10.04	124.63	118.60
53	B5	846	A	N1-C6-N6	10.04	124.62	118.60
53	B5	2628	A	N1-C6-N6	10.04	124.62	118.60
53	B5	338	A	N1-C6-N6	10.04	124.62	118.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	1204	A	N1-C6-N6	10.04	124.62	118.60
1	AA	180	A	N1-C6-N6	10.03	124.62	118.60
1	AA	755	A	N1-C6-N6	10.03	124.62	118.60
53	B5	2969	A	N1-C6-N6	10.03	124.62	118.60
1	AA	1467	A	N1-C6-N6	10.03	124.62	118.60
53	B5	39	A	N1-C6-N6	10.03	124.62	118.60
53	B5	3269	A	N1-C6-N6	10.03	124.62	118.60
1	AA	737	A	N1-C6-N6	10.03	124.61	118.60
1	AA	1424	A	N1-C6-N6	10.03	124.61	118.60
19	A7	44	A	C4-C5-C6	-10.03	111.99	117.00
53	B5	736	A	N1-C6-N6	10.02	124.61	118.60
1	AA	1200	A	C5-C6-N6	-10.02	115.69	123.70
1	AA	61	A	N1-C6-N6	10.02	124.61	118.60
53	B5	45	A	N1-C6-N6	10.01	124.61	118.60
1	AA	401	A	N1-C6-N6	10.01	124.61	118.60
1	AA	1669	A	N1-C6-N6	10.01	124.61	118.60
52	B4	70	A	N1-C6-N6	10.01	124.61	118.60
53	B5	389	A	N1-C6-N6	10.01	124.60	118.60
1	AA	1200	A	C2-N3-C4	10.00	115.60	110.60
1	AA	452	A	N1-C6-N6	10.00	124.60	118.60
53	B5	1102	A	N1-C6-N6	10.00	124.60	118.60
53	B5	1503	A	N1-C6-N6	10.00	124.60	118.60
53	B5	2601	A	N1-C6-N6	10.00	124.60	118.60
53	B5	2761	G	N1-C6-O6	10.00	125.90	119.90
19	A7	66	A	C6-C5-N7	10.00	139.30	132.30
52	B4	41	A	N1-C6-N6	10.00	124.60	118.60
53	B5	578	A	N1-C6-N6	10.00	124.60	118.60
53	B5	3000	A	N1-C6-N6	10.00	124.60	118.60
1	AA	1646	A	N1-C6-N6	9.99	124.60	118.60
53	B5	1259	A	N1-C6-N6	9.99	124.59	118.60
1	AA	1443	A	N1-C6-N6	9.99	124.59	118.60
53	B5	2252	A	N1-C6-N6	9.99	124.59	118.60
53	B5	1648	A	N1-C6-N6	9.99	124.59	118.60
53	B5	2967	A	N1-C6-N6	9.99	124.59	118.60
1	AA	906	A	N1-C6-N6	9.98	124.59	118.60
1	AA	938	A	N1-C6-N6	9.98	124.59	118.60
52	B4	97	A	N1-C6-N6	9.98	124.59	118.60
53	B5	1105	A	N1-C6-N6	9.98	124.59	118.60
53	B5	1337	A	N1-C6-N6	9.98	124.59	118.60
53	B5	1462	A	N1-C6-N6	9.98	124.59	118.60
1	AA	438	A	N1-C6-N6	9.98	124.59	118.60
1	AA	916	U	P-O3'-C3'	9.98	131.68	119.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1222	A	N1-C6-N6	9.98	124.59	118.60
53	B5	806	A	N1-C6-N6	9.98	124.59	118.60
1	AA	72	A	N1-C6-N6	9.97	124.58	118.60
53	B5	1883	A	N1-C6-N6	9.97	124.58	118.60
53	B5	2259	A	N1-C6-N6	9.97	124.58	118.60
53	B5	2373	A	N1-C6-N6	9.97	124.58	118.60
53	B5	2438	A	N1-C6-N6	9.97	124.58	118.60
53	B5	365	A	N1-C6-N6	9.97	124.58	118.60
53	B5	2131	A	N1-C6-N6	9.97	124.58	118.60
1	AA	1655	U	P-O3'-C3'	9.96	131.66	119.70
53	B5	61	A	N1-C6-N6	9.96	124.58	118.60
53	B5	355	A	N1-C6-N6	9.96	124.58	118.60
53	B5	2900	A	N1-C6-N6	9.96	124.58	118.60
1	AA	1038	A	N1-C6-N6	9.96	124.58	118.60
52	B4	9	A	N1-C6-N6	9.96	124.58	118.60
53	B5	2689	A	N1-C6-N6	9.96	124.58	118.60
53	B5	3183	A	N1-C6-N6	9.96	124.58	118.60
53	B5	325	A	N1-C6-N6	9.96	124.58	118.60
53	B5	344	A	N1-C6-N6	9.96	124.58	118.60
53	B5	2303	A	N1-C6-N6	9.96	124.57	118.60
1	AA	1062	A	N1-C6-N6	9.95	124.57	118.60
53	B5	1221	A	N1-C6-N6	9.95	124.57	118.60
53	B5	1731	A	N1-C6-N6	9.96	124.57	118.60
1	AA	511	A	N1-C6-N6	9.95	124.57	118.60
53	B5	735	A	N1-C6-N6	9.95	124.57	118.60
53	B5	1642	A	N1-C6-N6	9.95	124.57	118.60
53	B5	2790	A	N1-C6-N6	9.95	124.57	118.60
1	AA	1752	A	N1-C6-N6	9.95	124.57	118.60
53	B5	1558	A	N1-C6-N6	9.95	124.57	118.60
53	B5	2971	A	N1-C6-N6	9.95	124.57	118.60
53	B5	3029	A	N1-C6-N6	9.94	124.57	118.60
1	AA	1240	A	N1-C6-N6	9.94	124.57	118.60
53	B5	1593	A	N1-C6-N6	9.94	124.56	118.60
53	B5	1858	A	N1-C6-N6	9.94	124.56	118.60
1	AA	1388	A	N1-C6-N6	9.94	124.56	118.60
53	B5	657	A	N1-C6-N6	9.94	124.56	118.60
53	B5	1225	A	N1-C6-N6	9.94	124.56	118.60
53	B5	16	A	N1-C6-N6	9.94	124.56	118.60
1	AA	105	A	N1-C6-N6	9.93	124.56	118.60
1	AA	1326	A	N1-C6-N6	9.93	124.56	118.60
1	AA	1379	A	N1-C6-N6	9.93	124.56	118.60
53	B5	357	A	N1-C6-N6	9.93	124.56	118.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	2104	A	N1-C6-N6	9.93	124.56	118.60
53	B5	532	A	N1-C6-N6	9.93	124.56	118.60
53	B5	2694	A	N1-C6-N6	9.93	124.56	118.60
53	B5	374	A	N1-C6-N6	9.93	124.56	118.60
53	B5	1911	A	N1-C6-N6	9.92	124.56	118.60
1	AA	1687	A	N1-C6-N6	9.92	124.55	118.60
53	B5	2399	A	N1-C6-N6	9.92	124.55	118.60
53	B5	2401	A	N1-C6-N6	9.92	124.55	118.60
53	B5	2635	A	N1-C6-N6	9.92	124.55	118.60
1	AA	757	A	N1-C6-N6	9.91	124.55	118.60
53	B5	3165	A	N1-C6-N6	9.91	124.55	118.60
53	B5	1419	A	N1-C6-N6	9.91	124.55	118.60
53	B5	2462	A	N1-C6-N6	9.91	124.55	118.60
53	B5	3372	A	N1-C6-N6	9.91	124.55	118.60
1	AA	378	A	N1-C6-N6	9.91	124.55	118.60
53	B5	3243	A	N1-C6-N6	9.91	124.55	118.60
1	AA	1583	U	C4'-C3'-C2'	-9.91	92.69	102.60
1	AA	1728	A	N1-C6-N6	9.91	124.54	118.60
53	B5	62	A	N1-C6-N6	9.91	124.54	118.60
1	AA	1284	A	N1-C6-N6	9.90	124.54	118.60
53	B5	1534	A	N1-C6-N6	9.90	124.54	118.60
53	B5	2733	A	N1-C6-N6	9.90	124.54	118.60
53	B5	3134	A	N1-C6-N6	9.90	124.54	118.60
53	B5	666	A	N1-C6-N6	9.90	124.54	118.60
53	B5	3126	C	O4'-C1'-N1	9.90	116.12	108.20
1	AA	538	A	N1-C6-N6	9.90	124.54	118.60
1	AA	1408	A	N1-C6-N6	9.90	124.54	118.60
53	B5	933	A	N1-C6-N6	9.89	124.54	118.60
53	B5	2198	A	N1-C6-N6	9.89	124.54	118.60
53	B5	3011	A	N1-C6-N6	9.89	124.54	118.60
1	AA	85	A	N1-C6-N6	9.89	124.54	118.60
53	B5	2676	A	N1-C6-N6	9.89	124.54	118.60
1	AA	1148	A	N1-C6-N6	9.89	124.53	118.60
1	AA	1719	A	N1-C6-N6	9.89	124.53	118.60
53	B5	2219	A	N1-C6-N6	9.89	124.53	118.60
53	B5	998	A	N1-C6-N6	9.89	124.53	118.60
53	B5	2982	A	N1-C6-N6	9.89	124.53	118.60
53	B5	397	A	N1-C6-N6	9.88	124.53	118.60
53	B5	742	G	P-O3'-C3'	9.88	131.56	119.70
53	B5	1454	A	N1-C6-N6	9.88	124.53	118.60
19	A7	38	A	C4-C5-C6	-9.88	112.06	117.00
1	AA	254	A	N1-C6-N6	9.88	124.53	118.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1113	A	N1-C6-N6	9.88	124.53	118.60
53	B5	284	A	N1-C6-N6	9.88	124.53	118.60
53	B5	307	A	N1-C6-N6	9.88	124.53	118.60
1	AA	515	A	N1-C6-N6	9.88	124.53	118.60
53	B5	313	A	N1-C6-N6	9.88	124.53	118.60
53	B5	915	A	N1-C6-N6	9.88	124.53	118.60
1	AA	28	A	N1-C6-N6	9.88	124.53	118.60
53	B5	235	A	N1-C6-N6	9.87	124.52	118.60
53	B5	1401	A	N1-C6-N6	9.87	124.52	118.60
53	B5	1535	A	N1-C6-N6	9.87	124.52	118.60
53	B5	1841	A	N1-C6-N6	9.87	124.52	118.60
1	AA	367	A	N1-C6-N6	9.87	124.52	118.60
1	AA	788	A	N1-C6-N6	9.87	124.52	118.60
53	B5	1932	A	N1-C6-N6	9.87	124.52	118.60
53	B5	2792	A	N1-C6-N6	9.87	124.52	118.60
1	AA	300	A	N1-C6-N6	9.87	124.52	118.60
1	AA	352	A	N1-C6-N6	9.86	124.52	118.60
53	B5	1040	A	N1-C6-N6	9.86	124.52	118.60
53	B5	2144	A	N1-C6-N6	9.86	124.52	118.60
53	B5	3185	A	N1-C6-N6	9.86	124.52	118.60
1	AA	770	A	N1-C6-N6	9.86	124.52	118.60
52	B4	57	C	P-O3'-C3'	9.86	131.53	119.70
53	B5	523	A	N1-C6-N6	9.86	124.52	118.60
53	B5	660	A	N1-C6-N6	9.86	124.52	118.60
53	B5	1699	A	N1-C6-N6	9.86	124.52	118.60
53	B5	3048	A	N1-C6-N6	9.86	124.52	118.60
1	AA	164	A	N1-C6-N6	9.86	124.52	118.60
1	AA	1738	A	N1-C6-N6	9.86	124.52	118.60
53	B5	3391	A	N1-C6-N6	9.86	124.52	118.60
53	B5	6	A	N1-C6-N6	9.86	124.51	118.60
1	AA	366	A	N1-C6-N6	9.85	124.51	118.60
1	AA	904	A	N1-C6-N6	9.85	124.51	118.60
53	B5	882	A	N1-C6-N6	9.85	124.51	118.60
53	B5	3142	A	N1-C6-N6	9.85	124.51	118.60
53	B5	3017	A	N1-C6-N6	9.85	124.51	118.60
19	A7	76	A	C6-C5-N7	9.85	139.19	132.30
51	B3	77	A	N1-C6-N6	9.85	124.51	118.60
53	B5	1055	A	N1-C6-N6	9.85	124.51	118.60
1	AA	156	A	N1-C6-N6	9.85	124.51	118.60
1	AA	1717	A	N1-C6-N6	9.85	124.51	118.60
53	B5	2799	A	N1-C6-N6	9.85	124.51	118.60
1	AA	471	A	N1-C6-N6	9.84	124.51	118.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	1945	A	N1-C6-N6	9.84	124.50	118.60
1	AA	850	A	N1-C6-N6	9.84	124.50	118.60
53	B5	557	A	N1-C6-N6	9.84	124.50	118.60
1	AA	789	A	N1-C6-N6	9.84	124.50	118.60
53	B5	2147	A	N1-C6-N6	9.83	124.50	118.60
1	AA	138	A	N1-C6-N6	9.83	124.50	118.60
1	AA	359	A	N1-C6-N6	9.83	124.50	118.60
53	B5	3035	A	N1-C6-N6	9.83	124.50	118.60
53	B5	1235	U	P-O3'-C3'	9.83	131.50	119.70
4	AD	8	TYR	CB-CG-CD2	9.82	126.89	121.00
53	B5	1490	A	N1-C6-N6	9.82	124.50	118.60
53	B5	2166	A	N1-C6-N6	9.82	124.49	118.60
1	AA	1163	A	N1-C6-N6	9.82	124.49	118.60
53	B5	367	A	N1-C6-N6	9.82	124.49	118.60
1	AA	387	A	N1-C6-N6	9.82	124.49	118.60
1	AA	541	A	N1-C6-N6	9.81	124.49	118.60
1	AA	1544	G	O5'-P-OP1	9.81	122.48	110.70
53	B5	144	A	N1-C6-N6	9.81	124.49	118.60
53	B5	551	A	N1-C6-N6	9.81	124.49	118.60
1	AA	140	A	N1-C6-N6	9.81	124.49	118.60
53	B5	3175	C	O4'-C1'-N1	9.81	116.05	108.20
1	AA	542	A	N1-C6-N6	9.81	124.48	118.60
1	AA	1699	A	N1-C6-N6	9.81	124.48	118.60
19	A7	69	U	C5-C4-O4	-9.81	120.02	125.90
53	B5	598	A	N1-C6-N6	9.81	124.48	118.60
1	AA	78	A	N1-C6-N6	9.81	124.48	118.60
1	AA	397	A	N1-C6-N6	9.80	124.48	118.60
1	AA	922	A	N1-C6-N6	9.80	124.48	118.60
1	AA	1309	A	N1-C6-N6	9.80	124.48	118.60
53	B5	1842	A	N1-C6-N6	9.80	124.48	118.60
53	B5	1594	A	N1-C6-N6	9.80	124.48	118.60
53	B5	2721	A	N1-C6-N6	9.80	124.48	118.60
1	AA	1567	A	N1-C6-N6	9.80	124.48	118.60
1	AA	215	A	N1-C6-N6	9.80	124.48	118.60
19	A7	11	C	C6-N1-C2	-9.80	116.38	120.30
19	A7	53	G	N7-C8-N9	9.80	118.00	113.10
51	B3	94	A	N1-C6-N6	9.80	124.48	118.60
52	B4	66	A	N1-C6-N6	9.80	124.48	118.60
53	B5	1506	A	N1-C6-N6	9.80	124.48	118.60
53	B5	2520	A	N1-C6-N6	9.80	124.48	118.60
1	AA	119	A	N1-C6-N6	9.80	124.48	118.60
19	A7	6	U	C4-C5-C6	9.80	125.58	119.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	123	A	N1-C6-N6	9.80	124.48	118.60
1	AA	506	A	N1-C6-N6	9.79	124.48	118.60
51	B3	17	A	N1-C6-N6	9.79	124.48	118.60
52	B4	33	A	N1-C6-N6	9.79	124.48	118.60
53	B5	211	A	N1-C6-N6	9.80	124.48	118.60
53	B5	317	A	N1-C6-N6	9.79	124.48	118.60
53	B5	2643	A	N1-C6-N6	9.79	124.48	118.60
53	B5	3123	A	N1-C6-N6	9.79	124.48	118.60
1	AA	803	A	N1-C6-N6	9.79	124.47	118.60
52	B4	3	A	N1-C6-N6	9.79	124.47	118.60
53	B5	187	A	N1-C6-N6	9.79	124.47	118.60
53	B5	2561	A	N1-C6-N6	9.79	124.47	118.60
1	AA	220	A	N1-C6-N6	9.79	124.47	118.60
53	B5	1580	A	N1-C6-N6	9.79	124.47	118.60
1	AA	1397	A	N1-C6-N6	9.79	124.47	118.60
53	B5	2734	A	N1-C6-N6	9.79	124.47	118.60
53	B5	2872	A	N1-C6-N6	9.79	124.47	118.60
53	B5	1084	A	N1-C6-N6	9.79	124.47	118.60
1	AA	181	A	N1-C6-N6	9.79	124.47	118.60
53	B5	2106	A	N1-C6-N6	9.79	124.47	118.60
53	B5	1054	A	N1-C6-N6	9.78	124.47	118.60
53	B5	1867	A	N1-C6-N6	9.78	124.47	118.60
1	AA	169	A	N1-C6-N6	9.78	124.47	118.60
1	AA	271	A	N1-C6-N6	9.78	124.47	118.60
1	AA	1345	A	N1-C6-N6	9.78	124.47	118.60
53	B5	3094	A	N1-C6-N6	9.78	124.47	118.60
1	AA	570	A	N1-C6-N6	9.78	124.47	118.60
1	AA	1433	A	N1-C6-N6	9.78	124.47	118.60
52	B4	54	A	N1-C6-N6	9.78	124.47	118.60
53	B5	423	A	N1-C6-N6	9.78	124.47	118.60
53	B5	327	A	N1-C6-N6	9.78	124.47	118.60
53	B5	594	A	N1-C6-N6	9.78	124.47	118.60
53	B5	285	A	N1-C6-N6	9.78	124.47	118.60
1	AA	1190	A	N1-C6-N6	9.77	124.46	118.60
1	AA	1215	A	N1-C6-N6	9.77	124.46	118.60
1	AA	1220	A	N1-C6-N6	9.77	124.46	118.60
53	B5	255	A	N1-C6-N6	9.77	124.46	118.60
53	B5	1872	C	O4'-C1'-N1	9.77	116.02	108.20
53	B5	680	G	N1-C6-O6	9.77	125.76	119.90
53	B5	2397	A	N1-C6-N6	9.77	124.46	118.60
53	B5	2500	A	N1-C6-N6	9.77	124.46	118.60
52	B4	13	A	N1-C6-N6	9.77	124.46	118.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	936	A	N1-C6-N6	9.77	124.46	118.60
53	B5	1205	A	N1-C6-N6	9.77	124.46	118.60
53	B5	2535	A	N1-C6-N6	9.77	124.46	118.60
53	B5	2673	A	N1-C6-N6	9.77	124.46	118.60
53	B5	3039	C	O4'-C1'-N1	9.77	116.02	108.20
53	B5	1180	A	N1-C6-N6	9.77	124.46	118.60
53	B5	2991	A	N1-C6-N6	9.77	124.46	118.60
53	B5	3370	A	N1-C6-N6	9.76	124.46	118.60
1	AA	1058	A	N1-C6-N6	9.76	124.46	118.60
53	B5	3012	A	N1-C6-N6	9.76	124.46	118.60
53	B5	876	A	N1-C6-N6	9.76	124.45	118.60
53	B5	361	A	N1-C6-N6	9.76	124.45	118.60
1	AA	1333	A	N1-C6-N6	9.75	124.45	118.60
1	AA	1779	A	N1-C6-N6	9.75	124.45	118.60
53	B5	3223	A	N1-C6-N6	9.75	124.45	118.60
1	AA	713	A	N1-C6-N6	9.75	124.45	118.60
19	A7	7	U	C4-C5-C6	9.75	125.55	119.70
53	B5	501	A	N1-C6-N6	9.75	124.45	118.60
53	B5	569	A	N1-C6-N6	9.75	124.45	118.60
53	B5	697	A	N1-C6-N6	9.75	124.45	118.60
1	AA	295	A	N1-C6-N6	9.75	124.45	118.60
1	AA	534	A	N1-C6-N6	9.75	124.45	118.60
1	AA	684	A	N1-C6-N6	9.75	124.45	118.60
1	AA	1223	A	N1-C6-N6	9.75	124.45	118.60
53	B5	1524	A	N1-C6-N6	9.74	124.45	118.60
1	AA	1121	A	N1-C6-N6	9.74	124.44	118.60
51	B3	82	A	N1-C6-N6	9.74	124.45	118.60
53	B5	1200	A	N1-C6-N6	9.74	124.44	118.60
53	B5	720	A	N1-C6-N6	9.74	124.44	118.60
53	B5	2243	A	N1-C6-N6	9.74	124.44	118.60
1	AA	265	A	N1-C6-N6	9.74	124.44	118.60
1	AA	420	A	N1-C6-N6	9.74	124.44	118.60
53	B5	580	C	O4'-C1'-N1	9.74	115.99	108.20
53	B5	1303	A	N1-C6-N6	9.74	124.44	118.60
53	B5	1760	A	N1-C6-N6	9.74	124.44	118.60
53	B5	888	A	N1-C6-N6	9.73	124.44	118.60
53	B5	1135	A	N1-C6-N6	9.73	124.44	118.60
53	B5	2813	A	N1-C6-N6	9.73	124.44	118.60
1	AA	145	A	N1-C6-N6	9.73	124.44	118.60
1	AA	527	A	N1-C6-N6	9.73	124.44	118.60
53	B5	706	A	N1-C6-N6	9.73	124.44	118.60
53	B5	1202	A	N1-C6-N6	9.73	124.44	118.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	2330	C	O4'-C1'-N1	9.73	115.99	108.20
53	B5	755	A	N1-C6-N6	9.73	124.44	118.60
53	B5	786	A	N1-C6-N6	9.73	124.44	118.60
53	B5	1798	A	N1-C6-N6	9.73	124.44	118.60
53	B5	122	A	N1-C6-N6	9.73	124.44	118.60
1	AA	477	A	N1-C6-N6	9.73	124.44	118.60
1	AA	972	A	N1-C6-N6	9.73	124.44	118.60
53	B5	1850	A	N1-C6-N6	9.73	124.44	118.60
53	B5	1936	A	N1-C6-N6	9.73	124.44	118.60
53	B5	3206	A	N1-C6-N6	9.73	124.44	118.60
1	AA	1612	A	N1-C6-N6	9.73	124.44	118.60
53	B5	113	C	O4'-C1'-N1	9.73	115.98	108.20
53	B5	550	A	N1-C6-N6	9.73	124.44	118.60
53	B5	738	A	N1-C6-N6	9.73	124.44	118.60
53	B5	1169	A	N1-C6-N6	9.73	124.44	118.60
1	AA	92	A	N1-C6-N6	9.72	124.43	118.60
1	AA	1090	A	N1-C6-N6	9.72	124.44	118.60
1	AA	407	A	N1-C6-N6	9.72	124.43	118.60
51	B3	45	A	N1-C6-N6	9.72	124.44	118.60
53	B5	665	A	N1-C6-N6	9.72	124.43	118.60
53	B5	1654	A	N1-C6-N6	9.72	124.43	118.60
1	AA	1780	A	N1-C6-N6	9.72	124.43	118.60
53	B5	1006	A	N1-C6-N6	9.72	124.43	118.60
53	B5	1223	A	N1-C6-N6	9.72	124.43	118.60
53	B5	2637	A	N1-C6-N6	9.72	124.43	118.60
52	B4	77	A	N1-C6-N6	9.72	124.43	118.60
53	B5	2642	A	N1-C6-N6	9.71	124.43	118.60
53	B5	2896	A	N1-C6-N6	9.72	124.43	118.60
53	B5	3344	A	N1-C6-N6	9.72	124.43	118.60
53	B5	3113	A	N1-C6-N6	9.71	124.43	118.60
1	AA	108	A	N1-C6-N6	9.71	124.43	118.60
53	B5	715	A	N1-C6-N6	9.71	124.43	118.60
1	AA	221	A	N1-C6-N6	9.71	124.43	118.60
1	AA	678	A	N1-C6-N6	9.71	124.43	118.60
53	B5	1461	A	N1-C6-N6	9.71	124.43	118.60
53	B5	2424	A	N1-C6-N6	9.71	124.42	118.60
1	AA	47	A	N1-C6-N6	9.71	124.42	118.60
19	A7	61	C	N1-C2-O2	9.71	124.72	118.90
52	B4	12	A	N1-C6-N6	9.71	124.42	118.60
53	B5	334	A	N1-C6-N6	9.70	124.42	118.60
1	AA	410	A	N1-C6-N6	9.70	124.42	118.60
1	AA	860	U	O4'-C1'-N1	9.70	115.96	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1551	G	O5'-P-OP1	9.70	122.34	110.70
1	AA	1501	A	N1-C6-N6	9.70	124.42	118.60
53	B5	653	A	N1-C6-N6	9.70	124.42	118.60
53	B5	3375	A	N1-C6-N6	9.70	124.42	118.60
53	B5	2488	A	N1-C6-N6	9.69	124.42	118.60
53	B5	2746	A	N1-C6-N6	9.70	124.42	118.60
1	AA	416	A	N1-C6-N6	9.69	124.42	118.60
1	AA	1421	A	N1-C6-N6	9.69	124.42	118.60
1	AA	1789	A	N1-C6-N6	9.69	124.41	118.60
53	B5	3070	A	N1-C6-N6	9.69	124.41	118.60
1	AA	1089	A	N1-C6-N6	9.69	124.41	118.60
53	B5	165	A	N1-C6-N6	9.69	124.41	118.60
53	B5	692	A	N1-C6-N6	9.69	124.41	118.60
53	B5	1308	A	N1-C6-N6	9.69	124.41	118.60
53	B5	2740	A	N1-C6-N6	9.69	124.41	118.60
53	B5	896	A	N1-C6-N6	9.69	124.41	118.60
53	B5	990	A	N1-C6-N6	9.69	124.41	118.60
53	B5	1452	A	N1-C6-N6	9.69	124.41	118.60
1	AA	661	A	N1-C6-N6	9.68	124.41	118.60
53	B5	645	A	N1-C6-N6	9.68	124.41	118.60
1	AA	441	A	N1-C6-N6	9.68	124.41	118.60
1	AA	1088	A	N1-C6-N6	9.68	124.41	118.60
1	AA	1414	A	N1-C6-N6	9.68	124.41	118.60
1	AA	1490	A	N1-C6-N6	9.68	124.41	118.60
53	B5	2864	A	N1-C6-N6	9.68	124.41	118.60
53	B5	2152	A	N1-C6-N6	9.68	124.41	118.60
53	B5	621	A	N1-C6-N6	9.68	124.41	118.60
53	B5	2692	A	N1-C6-N6	9.68	124.41	118.60
51	B3	19	C	O4'-C1'-N1	9.67	115.94	108.20
53	B5	1212	A	N1-C6-N6	9.67	124.41	118.60
53	B5	1394	A	N1-C6-N6	9.67	124.40	118.60
53	B5	1638	A	N1-C6-N6	9.67	124.40	118.60
53	B5	529	A	N1-C6-N6	9.67	124.40	118.60
53	B5	2887	A	N1-C6-N6	9.67	124.40	118.60
53	B5	2890	A	N1-C6-N6	9.67	124.40	118.60
53	B5	1459	C	O4'-C1'-N1	9.67	115.93	108.20
1	AA	222	A	N1-C6-N6	9.66	124.40	118.60
1	AA	555	A	N1-C6-N6	9.66	124.40	118.60
1	AA	987	A	N1-C6-N6	9.66	124.40	118.60
53	B5	904	A	N1-C6-N6	9.66	124.40	118.60
53	B5	1498	A	N1-C6-N6	9.66	124.40	118.60
1	AA	1503	A	N1-C6-N6	9.66	124.40	118.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
19	A7	8	U	N3-C4-C5	-9.66	108.80	114.60
1	AA	1477	A	N1-C6-N6	9.66	124.39	118.60
53	B5	789	A	N1-C6-N6	9.66	124.39	118.60
1	AA	978	A	N1-C6-N6	9.65	124.39	118.60
1	AA	1609	A	N1-C6-N6	9.65	124.39	118.60
53	B5	250	U	P-O3'-C3'	9.65	131.28	119.70
53	B5	3359	A	N1-C6-N6	9.65	124.39	118.60
1	AA	1078	A	N1-C6-N6	9.65	124.39	118.60
53	B5	121	A	N1-C6-N6	9.65	124.39	118.60
1	AA	333	A	N1-C6-N6	9.65	124.39	118.60
1	AA	385	A	N1-C6-N6	9.65	124.39	118.60
1	AA	1517	U	O5'-P-OP1	9.65	122.28	110.70
53	B5	1170	A	N1-C6-N6	9.65	124.39	118.60
53	B5	1489	A	N1-C6-N6	9.65	124.39	118.60
53	B5	369	A	N1-C6-N6	9.65	124.39	118.60
1	AA	1083	A	N1-C6-N6	9.64	124.39	118.60
1	AA	1341	A	N1-C6-N6	9.64	124.39	118.60
52	B4	80	A	N1-C6-N6	9.64	124.39	118.60
53	B5	1009	A	N1-C6-N6	9.64	124.39	118.60
53	B5	1806	A	N1-C6-N6	9.64	124.39	118.60
53	B5	2220	A	N1-C6-N6	9.64	124.39	118.60
1	AA	740	A	N1-C6-N6	9.64	124.38	118.60
53	B5	2345	A	N1-C6-N6	9.64	124.39	118.60
53	B5	2119	A	N1-C6-N6	9.64	124.38	118.60
53	B5	3316	A	N1-C6-N6	9.64	124.38	118.60
1	AA	40	A	N1-C6-N6	9.64	124.38	118.60
1	AA	173	A	N1-C6-N6	9.64	124.38	118.60
1	AA	526	A	N1-C6-N6	9.64	124.38	118.60
1	AA	791	A	N1-C6-N6	9.64	124.38	118.60
53	B5	646	A	N1-C6-N6	9.64	124.38	118.60
1	AA	421	A	N1-C6-N6	9.64	124.38	118.60
53	B5	2207	A	N1-C6-N6	9.64	124.38	118.60
53	B5	2549	A	N1-C6-N6	9.64	124.38	118.60
53	B5	843	A	N1-C6-N6	9.64	124.38	118.60
1	AA	858	G	N3-C2-N2	9.63	126.64	119.90
53	B5	1823	A	N1-C6-N6	9.63	124.38	118.60
53	B5	521	A	N1-C6-N6	9.63	124.38	118.60
53	B5	2320	A	N1-C6-N6	9.63	124.38	118.60
53	B5	2987	A	N1-C6-N6	9.63	124.38	118.60
53	B5	970	A	N1-C6-N6	9.62	124.37	118.60
53	B5	3008	A	N1-C6-N6	9.62	124.37	118.60
53	B5	3040	A	N1-C6-N6	9.62	124.37	118.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1458	A	N1-C6-N6	9.62	124.37	118.60
52	B4	79	A	N1-C6-N6	9.62	124.37	118.60
53	B5	43	A	N1-C6-N6	9.62	124.37	118.60
53	B5	3106	A	N1-C6-N6	9.62	124.37	118.60
51	B3	42	A	N1-C6-N6	9.62	124.37	118.60
1	AA	1689	A	N1-C6-N6	9.61	124.37	118.60
1	AA	1763	A	N1-C6-N6	9.62	124.37	118.60
53	B5	34	A	N1-C6-N6	9.62	124.37	118.60
53	B5	3176	C	O4'-C1'-N1	9.61	115.89	108.20
1	AA	550	A	N1-C6-N6	9.61	124.37	118.60
53	B5	1474	A	N1-C6-N6	9.61	124.37	118.60
53	B5	1557	A	N1-C6-N6	9.61	124.37	118.60
53	B5	2138	A	N1-C6-N6	9.61	124.37	118.60
1	AA	525	A	N1-C6-N6	9.61	124.37	118.60
51	B3	23	A	N1-C6-N6	9.61	124.36	118.60
53	B5	1304	A	N1-C6-N6	9.61	124.36	118.60
53	B5	1048	A	O4'-C1'-N9	9.61	115.88	108.20
53	B5	2468	A	N1-C6-N6	9.61	124.36	118.60
1	AA	126	A	N1-C6-N6	9.60	124.36	118.60
1	AA	604	A	N1-C6-N6	9.60	124.36	118.60
19	A7	27	C	N1-C2-N3	9.60	125.92	119.20
1	AA	1381	A	N1-C6-N6	9.60	124.36	118.60
1	AA	1549	U	O4'-C1'-N1	9.60	115.88	108.20
53	B5	1030	A	N1-C6-N6	9.60	124.36	118.60
53	B5	2398	A	N1-C6-N6	9.60	124.36	118.60
53	B5	2695	A	N1-C6-N6	9.60	124.36	118.60
53	B5	3215	A	N1-C6-N6	9.60	124.36	118.60
1	AA	34	G	O4'-C1'-N9	9.60	115.88	108.20
1	AA	65	A	N1-C6-N6	9.60	124.36	118.60
53	B5	85	A	N1-C6-N6	9.60	124.36	118.60
53	B5	150	A	N1-C6-N6	9.60	124.36	118.60
53	B5	2130	G	C5-C6-O6	-9.60	122.84	128.60
53	B5	3146	G	O4'-C1'-N9	9.60	115.88	108.20
1	AA	594	A	N1-C6-N6	9.59	124.35	118.60
53	B5	1712	G	O4'-C1'-N9	9.59	115.87	108.20
1	AA	867	G	P-O3'-C3'	-9.59	108.19	119.70
1	AA	925	A	N1-C6-N6	9.59	124.35	118.60
1	AA	1368	A	N1-C6-N6	9.59	124.35	118.60
53	B5	2441	A	N1-C6-N6	9.59	124.35	118.60
53	B5	422	A	N1-C6-N6	9.58	124.35	118.60
53	B5	1908	A	N1-C6-N6	9.58	124.35	118.60
1	AA	251	A	N1-C6-N6	9.58	124.35	118.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	76	A	N1-C6-N6	9.58	124.35	118.60
53	B5	490	A	N1-C6-N6	9.58	124.35	118.60
53	B5	914	A	N1-C6-N6	9.58	124.35	118.60
53	B5	1143	A	N1-C6-N6	9.57	124.34	118.60
53	B5	2739	A	N1-C6-N6	9.57	124.34	118.60
1	AA	182	A	N1-C6-N6	9.57	124.34	118.60
1	AA	1774	A	N1-C6-N6	9.57	124.34	118.60
53	B5	2262	A	N1-C6-N6	9.57	124.34	118.60
53	B5	1182	A	N1-C6-N6	9.57	124.34	118.60
1	AA	344	A	N1-C6-N6	9.57	124.34	118.60
19	A7	23	A	C3'-C2'-C1'	9.57	109.15	101.50
53	B5	378	A	N1-C6-N6	9.57	124.34	118.60
53	B5	519	A	N1-C6-N6	9.57	124.34	118.60
53	B5	2271	A	N1-C6-N6	9.57	124.34	118.60
53	B5	2228	A	N1-C6-N6	9.56	124.34	118.60
53	B5	1587	A	N1-C6-N6	9.56	124.34	118.60
53	B5	2363	A	N1-C6-N6	9.56	124.34	118.60
1	AA	1691	A	N1-C6-N6	9.56	124.33	118.60
53	B5	296	A	N1-C6-N6	9.56	124.33	118.60
1	AA	217	A	N1-C6-N6	9.56	124.33	118.60
1	AA	620	A	N1-C6-N6	9.55	124.33	118.60
1	AA	754	A	N1-C6-N6	9.55	124.33	118.60
1	AA	997	A	N1-C6-N6	9.56	124.33	118.60
1	AA	1712	A	N1-C6-N6	9.56	124.33	118.60
1	AA	864	U	C4'-C3'-C2'	9.55	112.15	102.60
53	B5	2519	A	N1-C6-N6	9.55	124.33	118.60
53	B5	376	G	N1-C6-O6	9.55	125.63	119.90
53	B5	1810	A	N1-C6-N6	9.55	124.33	118.60
52	B4	48	A	N1-C6-N6	9.55	124.33	118.60
1	AA	253	A	N1-C6-N6	9.55	124.33	118.60
1	AA	1629	A	N1-C6-N6	9.55	124.33	118.60
53	B5	1373	A	N1-C6-N6	9.55	124.33	118.60
53	B5	2223	A	N1-C6-N6	9.55	124.33	118.60
53	B5	3296	A	N1-C6-N6	9.55	124.33	118.60
53	B5	533	A	N1-C6-N6	9.54	124.33	118.60
53	B5	2802	A	N1-C6-N6	9.55	124.33	118.60
53	B5	342	A	N1-C6-N6	9.54	124.33	118.60
53	B5	1098	A	N1-C6-N6	9.54	124.32	118.60
53	B5	2738	A	N1-C6-N6	9.54	124.32	118.60
53	B5	3335	A	N1-C6-N6	9.54	124.32	118.60
53	B5	2486	A	N1-C6-N6	9.54	124.32	118.60
1	AA	1149	A	N1-C6-N6	9.53	124.32	118.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	2941	A	N1-C6-N6	9.53	124.32	118.60
52	B4	61	A	N1-C6-N6	9.53	124.32	118.60
53	B5	289	A	N1-C6-N6	9.53	124.32	118.60
53	B5	2232	A	N1-C6-N6	9.53	124.32	118.60
53	B5	2413	A	N1-C6-N6	9.53	124.32	118.60
53	B5	3021	A	N1-C6-N6	9.53	124.32	118.60
1	AA	923	A	N1-C6-N6	9.53	124.32	118.60
1	AA	817	A	N1-C6-N6	9.52	124.31	118.60
51	B3	100	A	N1-C6-N6	9.52	124.31	118.60
53	B5	1061	A	N1-C6-N6	9.52	124.31	118.60
1	AA	312	A	N1-C6-N6	9.52	124.31	118.60
1	AA	529	A	N1-C6-N6	9.52	124.31	118.60
1	AA	606	A	N1-C6-N6	9.52	124.31	118.60
1	AA	257	A	N1-C6-N6	9.52	124.31	118.60
1	AA	1658	A	N1-C6-N6	9.52	124.31	118.60
53	B5	2484	A	N1-C6-N6	9.51	124.31	118.60
1	AA	1207	A	N1-C6-N6	9.51	124.31	118.60
53	B5	2547	A	N1-C6-N6	9.51	124.31	118.60
53	B5	2580	A	N1-C6-N6	9.51	124.31	118.60
1	AA	982	A	N1-C6-N6	9.51	124.30	118.60
19	A7	5	A	C5-C6-N1	9.51	122.45	117.70
53	B5	1467	A	N1-C6-N6	9.51	124.30	118.60
53	B5	3276	A	N1-C6-N6	9.51	124.30	118.60
53	B5	11	A	N1-C6-N6	9.50	124.30	118.60
53	B5	273	A	N1-C6-N6	9.50	124.30	118.60
53	B5	2803	A	N1-C6-N6	9.50	124.30	118.60
53	B5	1026	A	N1-C6-N6	9.50	124.30	118.60
1	AA	288	A	N1-C6-N6	9.50	124.30	118.60
1	AA	994	A	N1-C6-N6	9.50	124.30	118.60
1	AA	1657	A	N1-C6-N6	9.50	124.30	118.60
53	B5	1896	A	N1-C6-N6	9.49	124.30	118.60
1	AA	1549	U	C3'-C2'-C1'	-9.49	93.91	101.50
52	B4	138	A	N1-C6-N6	9.49	124.29	118.60
53	B5	107	A	N1-C6-N6	9.49	124.29	118.60
53	B5	1330	A	N1-C6-N6	9.49	124.29	118.60
53	B5	1465	A	N1-C6-N6	9.49	124.29	118.60
1	AA	1130	A	N1-C6-N6	9.49	124.29	118.60
1	AA	1342	A	N1-C6-N6	9.49	124.29	118.60
53	B5	2743	A	N1-C6-N6	9.49	124.29	118.60
53	B5	1407	A	N1-C6-N6	9.48	124.29	118.60
1	AA	1297	A	N1-C6-N6	9.48	124.29	118.60
53	B5	3114	A	N1-C6-N6	9.48	124.29	118.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	425	A	N1-C6-N6	9.48	124.29	118.60
1	AA	556	A	N1-C6-N6	9.48	124.29	118.60
1	AA	991	A	N1-C6-N6	9.48	124.29	118.60
1	AA	706	A	N1-C6-N6	9.48	124.29	118.60
1	AA	811	A	N1-C6-N6	9.48	124.29	118.60
1	AA	1136	A	N1-C6-N6	9.48	124.29	118.60
19	A7	14	A	N1-C6-N6	-9.48	112.91	118.60
51	B3	89	A	N1-C6-N6	9.48	124.29	118.60
53	B5	570	A	N1-C6-N6	9.48	124.28	118.60
1	AA	84	A	N1-C6-N6	9.47	124.28	118.60
1	AA	866	G	C4'-C3'-C2'	9.47	112.07	102.60
53	B5	2485	A	N1-C6-N6	9.47	124.28	118.60
53	B5	2820	A	N1-C6-N6	9.47	124.28	118.60
1	AA	622	A	N1-C6-N6	9.47	124.28	118.60
19	A7	1	G	C5-C6-N1	9.47	116.23	111.50
1	AA	512	A	N1-C6-N6	9.47	124.28	118.60
53	B5	1895	A	N1-C6-N6	9.47	124.28	118.60
53	B5	2595	A	N1-C6-N6	9.46	124.28	118.60
1	AA	733	A	N1-C6-N6	9.46	124.28	118.60
1	AA	1520	U	P-O5'-C5'	-9.46	105.76	120.90
52	B4	37	A	N1-C6-N6	9.46	124.28	118.60
1	AA	847	A	N1-C6-N6	9.46	124.28	118.60
53	B5	2430	A	N1-C6-N6	9.46	124.28	118.60
1	AA	323	A	N1-C6-N6	9.46	124.28	118.60
1	AA	630	A	N1-C6-N6	9.46	124.28	118.60
1	AA	1679	A	N1-C6-N6	9.46	124.28	118.60
53	B5	1324	U	O4'-C1'-N1	9.46	115.77	108.20
53	B5	23	A	N1-C6-N6	9.46	124.27	118.60
53	B5	2529	A	N1-C6-N6	9.46	124.27	118.60
1	AA	206	A	N1-C6-N6	9.46	124.27	118.60
53	B5	603	A	N1-C6-N6	9.46	124.27	118.60
1	AA	774	A	N1-C6-N6	9.45	124.27	118.60
1	AA	1514	A	N1-C6-N6	9.46	124.27	118.60
53	B5	1468	A	N1-C6-N6	9.46	124.27	118.60
53	B5	3045	G	P-O3'-C3'	9.46	131.05	119.70
1	AA	1073	A	N1-C6-N6	9.45	124.27	118.60
51	B3	75	G	N1-C6-O6	9.45	125.57	119.90
53	B5	3128	G	N1-C6-O6	9.45	125.57	119.90
1	AA	780	A	N1-C6-N6	9.45	124.27	118.60
53	B5	3163	A	N1-C6-N6	9.45	124.27	118.60
52	B4	111	A	N1-C6-N6	9.45	124.27	118.60
1	AA	752	A	N1-C6-N6	9.44	124.27	118.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1025	A	N1-C6-N6	9.44	124.27	118.60
1	AA	1543	A	N1-C6-N6	9.45	124.27	118.60
53	B5	204	A	N1-C6-N6	9.44	124.27	118.60
53	B5	1294	A	N1-C6-N6	9.44	124.27	118.60
53	B5	2275	A	N1-C6-N6	9.45	124.27	118.60
1	AA	1128	A	N1-C6-N6	9.44	124.27	118.60
1	AA	1403	A	N1-C6-N6	9.44	124.26	118.60
19	A7	64	A	C6-C5-N7	9.44	138.91	132.30
51	B3	35	C	O4'-C1'-N1	9.44	115.75	108.20
52	B4	2	A	N1-C6-N6	9.44	124.27	118.60
53	B5	672	A	N1-C6-N6	9.44	124.27	118.60
53	B5	2183	A	N1-C6-N6	9.44	124.26	118.60
53	B5	2280	A	N1-C6-N6	9.44	124.27	118.60
1	AA	1481	A	N1-C6-N6	9.44	124.26	118.60
1	AA	71	A	N1-C6-N6	9.44	124.26	118.60
53	B5	2994	A	N1-C6-N6	9.44	124.26	118.60
53	B5	1460	A	N1-C6-N6	9.43	124.26	118.60
1	AA	932	A	N1-C6-N6	9.43	124.26	118.60
1	AA	1729	A	N1-C6-N6	9.43	124.26	118.60
53	B5	164	A	N1-C6-N6	9.43	124.26	118.60
1	AA	328	A	N1-C6-N6	9.43	124.26	118.60
1	AA	1742	A	N1-C6-N6	9.43	124.26	118.60
53	B5	1002	A	N1-C6-N6	9.43	124.26	118.60
53	B5	1714	A	N1-C6-N6	9.43	124.26	118.60
52	B4	17	A	N1-C6-N6	9.43	124.26	118.60
53	B5	472	A	N1-C6-N6	9.43	124.26	118.60
53	B5	1900	A	N1-C6-N6	9.43	124.25	118.60
1	AA	475	A	N1-C6-N6	9.42	124.25	118.60
53	B5	51	A	N1-C6-N6	9.42	124.25	118.60
1	AA	1035	A	N1-C6-N6	9.42	124.25	118.60
51	B3	114	A	N1-C6-N6	9.42	124.25	118.60
53	B5	1280	C	O4'-C1'-N1	9.42	115.73	108.20
53	B5	3024	A	N1-C6-N6	9.42	124.25	118.60
53	B5	1656	A	N1-C6-N6	9.41	124.25	118.60
1	AA	11	A	N1-C6-N6	9.41	124.25	118.60
1	AA	483	A	N1-C6-N6	9.41	124.25	118.60
53	B5	2697	A	N1-C6-N6	9.41	124.25	118.60
53	B5	589	A	N1-C6-N6	9.41	124.25	118.60
1	AA	478	A	N1-C6-N6	9.41	124.25	118.60
53	B5	1637	A	N1-C6-N6	9.41	124.25	118.60
53	B5	2524	A	N1-C6-N6	9.41	124.25	118.60
1	AA	814	A	N1-C6-N6	9.41	124.24	118.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	887	A	N1-C6-N6	9.41	124.25	118.60
53	B5	527	A	N1-C6-N6	9.40	124.24	118.60
1	AA	46	A	N1-C6-N6	9.40	124.24	118.60
53	B5	751	A	N1-C6-N6	9.40	124.24	118.60
53	B5	49	A	N1-C6-N6	9.40	124.24	118.60
53	B5	3077	A	N1-C6-N6	9.40	124.24	118.60
1	AA	266	A	N1-C6-N6	9.40	124.24	118.60
53	B5	1515	A	N1-C6-N6	9.40	124.24	118.60
53	B5	2621	G	N1-C6-O6	9.40	125.54	119.90
53	B5	2785	A	N1-C6-N6	9.40	124.24	118.60
1	AA	446	A	N1-C6-N6	9.39	124.24	118.60
1	AA	544	A	N1-C6-N6	9.39	124.24	118.60
1	AA	619	A	N1-C6-N6	9.39	124.24	118.60
1	AA	1316	A	N1-C6-N6	9.39	124.24	118.60
53	B5	115	A	N1-C6-N6	9.39	124.24	118.60
53	B5	2125	A	N1-C6-N6	9.39	124.24	118.60
53	B5	3271	A	N1-C6-N6	9.39	124.24	118.60
1	AA	1598	A	N1-C6-N6	9.39	124.23	118.60
1	AA	1059	A	N1-C6-N6	9.39	124.23	118.60
53	B5	1179	A	N1-C6-N6	9.39	124.23	118.60
53	B5	1809	A	N1-C6-N6	9.39	124.23	118.60
52	B4	1	A	N1-C6-N6	9.38	124.23	118.60
53	B5	251	G	P-O3'-C3'	9.38	130.96	119.70
53	B5	1723	A	N1-C6-N6	9.38	124.23	118.60
53	B5	1749	A	N1-C6-N6	9.38	124.23	118.60
1	AA	240	U	O4'-C1'-N1	9.38	115.70	108.20
51	B3	27	A	N1-C6-N6	9.38	124.23	118.60
53	B5	106	A	N1-C6-N6	9.38	124.23	118.60
53	B5	1326	A	N1-C6-N6	9.38	124.23	118.60
53	B5	2324	A	N1-C6-N6	9.38	124.23	118.60
53	B5	2383	C	O4'-C1'-N1	9.38	115.70	108.20
1	AA	198	A	N1-C6-N6	9.38	124.22	118.60
1	AA	197	A	N1-C6-N6	9.37	124.22	118.60
53	B5	3110	C	O4'-C1'-N1	9.38	115.70	108.20
1	AA	210	A	N1-C6-N6	9.37	124.22	118.60
53	B5	2591	A	N1-C6-N6	9.37	124.22	118.60
19	A7	64	A	C4'-C3'-C2'	-9.37	93.23	102.60
53	B5	1136	A	N1-C6-N6	9.37	124.22	118.60
53	B5	1238	C	O4'-C1'-N1	9.37	115.69	108.20
53	B5	2801	A	N1-C6-N6	9.37	124.22	118.60
53	B5	2164	A	N1-C6-N6	9.37	124.22	118.60
53	B5	3032	A	N1-C6-N6	9.37	124.22	118.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	124	A	N1-C6-N6	9.37	124.22	118.60
1	AA	1571	A	N1-C6-N6	9.36	124.22	118.60
53	B5	691	A	N1-C6-N6	9.37	124.22	118.60
53	B5	841	A	N1-C6-N6	9.37	124.22	118.60
53	B5	3226	A	N1-C6-N6	9.37	124.22	118.60
1	AA	200	A	N1-C6-N6	9.36	124.22	118.60
1	AA	599	A	N1-C6-N6	9.36	124.22	118.60
1	AA	899	A	N1-C6-N6	9.36	124.22	118.60
1	AA	218	A	N1-C6-N6	9.36	124.22	118.60
1	AA	579	A	N1-C6-N6	9.36	124.21	118.60
19	A7	66	A	C4'-C3'-C2'	-9.36	93.24	102.60
53	B5	428	A	N1-C6-N6	9.36	124.21	118.60
1	AA	1180	A	N1-C6-N6	9.36	124.21	118.60
53	B5	1816	A	N1-C6-N6	9.36	124.21	118.60
1	AA	585	A	N1-C6-N6	9.35	124.21	118.60
1	AA	1040	A	N1-C6-N6	9.35	124.21	118.60
53	B5	174	C	O4'-C1'-N1	9.35	115.68	108.20
53	B5	1251	A	N1-C6-N6	9.35	124.21	118.60
53	B5	2647	A	N1-C6-N6	9.35	124.21	118.60
53	B5	391	A	N1-C6-N6	9.35	124.21	118.60
53	B5	2769	A	N1-C6-N6	9.35	124.21	118.60
53	B5	3330	A	N1-C6-N6	9.35	124.21	118.60
1	AA	19	A	N1-C6-N6	9.35	124.21	118.60
1	AA	213	A	N1-C6-N6	9.35	124.21	118.60
53	B5	883	A	N1-C6-N6	9.35	124.21	118.60
53	B5	1302	A	N1-C6-N6	9.35	124.21	118.60
53	B5	1256	G	N1-C6-O6	9.35	125.51	119.90
53	B5	504	A	N1-C6-N6	9.35	124.21	118.60
53	B5	1036	A	N1-C6-N6	9.35	124.21	118.60
53	B5	1270	A	N1-C6-N6	9.35	124.21	118.60
1	AA	807	A	N1-C6-N6	9.35	124.21	118.60
53	B5	2214	A	N1-C6-N6	9.35	124.21	118.60
53	B5	2808	A	N1-C6-N6	9.35	124.21	118.60
1	AA	225	A	N1-C6-N6	9.34	124.21	118.60
1	AA	1726	A	N1-C6-N6	9.34	124.21	118.60
53	B5	3339	A	N1-C6-N6	9.34	124.20	118.60
1	AA	1513	A	N1-C6-N6	9.34	124.20	118.60
53	B5	399	A	N1-C6-N6	9.34	124.20	118.60
53	B5	747	A	N1-C6-N6	9.34	124.20	118.60
1	AA	226	A	N1-C6-N6	9.33	124.20	118.60
1	AA	977	A	N1-C6-N6	9.33	124.20	118.60
52	B4	68	G	P-O3'-C3'	9.33	130.90	119.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	1475	A	N1-C6-N6	9.33	124.20	118.60
1	AA	1751	A	N1-C6-N6	9.33	124.20	118.60
52	B4	88	A	N1-C6-N6	9.33	124.20	118.60
1	AA	673	A	N1-C6-N6	9.33	124.20	118.60
13	AN	16	LYS	CB-CA-C	-9.33	91.74	110.40
53	B5	1428	A	N1-C6-N6	9.33	124.20	118.60
1	AA	760	A	N1-C6-N6	9.33	124.20	118.60
53	B5	630	A	N1-C6-N6	9.33	124.20	118.60
53	B5	649	A	N1-C6-N6	9.33	124.20	118.60
53	B5	2936	A	N1-C6-N6	9.33	124.20	118.60
53	B5	3028	G	N1-C6-O6	9.33	125.50	119.90
53	B5	1884	A	N1-C6-N6	9.32	124.19	118.60
1	AA	1	U	O4'-C1'-N1	9.32	115.66	108.20
1	AA	792	U	OP1-P-OP2	-9.32	105.62	119.60
51	B3	33	U	O4'-C1'-N1	9.32	115.66	108.20
53	B5	57	A	N1-C6-N6	9.32	124.19	118.60
53	B5	2270	A	N1-C6-N6	9.32	124.19	118.60
53	B5	3270	A	N1-C6-N6	9.32	124.19	118.60
53	B5	3292	A	N1-C6-N6	9.32	124.19	118.60
53	B5	1752	A	N1-C6-N6	9.31	124.19	118.60
1	AA	812	A	N1-C6-N6	9.30	124.18	118.60
1	AA	1521	G	C6-C5-N7	-9.31	124.82	130.40
53	B5	3244	A	N1-C6-N6	9.30	124.18	118.60
19	A7	41	U	C2-N3-C4	-9.30	121.42	127.00
53	B5	1453	A	N1-C6-N6	9.30	124.18	118.60
53	B5	1814	A	N1-C6-N6	9.30	124.18	118.60
53	B5	478	A	N1-C6-N6	9.29	124.18	118.60
53	B5	1154	A	N1-C6-N6	9.29	124.18	118.60
1	AA	492	A	N1-C6-N6	9.29	124.18	118.60
53	B5	1424	C	O4'-C1'-N1	9.29	115.63	108.20
1	AA	112	A	N1-C6-N6	9.29	124.17	118.60
1	AA	1012	A	N1-C6-N6	9.29	124.17	118.60
53	B5	516	A	N1-C6-N6	9.29	124.17	118.60
53	B5	656	A	N1-C6-N6	9.29	124.17	118.60
53	B5	780	A	N1-C6-N6	9.29	124.17	118.60
1	AA	869	A	N1-C6-N6	9.28	124.17	118.60
53	B5	349	A	N1-C6-N6	9.28	124.17	118.60
53	B5	2352	A	N1-C6-N6	9.28	124.17	118.60
53	B5	2863	G	O4'-C1'-N9	9.28	115.63	108.20
53	B5	2995	A	N1-C6-N6	9.29	124.17	118.60
1	AA	950	A	N1-C6-N6	9.28	124.17	118.60
1	AA	806	A	N1-C6-N6	9.28	124.17	118.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1232	A	N1-C6-N6	9.28	124.17	118.60
53	B5	1919	G	N1-C6-O6	9.28	125.47	119.90
53	B5	771	A	N1-C6-N6	9.28	124.17	118.60
53	B5	1046	A	N1-C6-N6	9.28	124.17	118.60
53	B5	1244	A	N1-C6-N6	9.28	124.17	118.60
53	B5	1571	A	N1-C6-N6	9.28	124.17	118.60
53	B5	2231	C	C6-N1-C1'	-9.28	109.67	120.80
1	AA	428	A	N1-C6-N6	9.28	124.17	118.60
1	AA	753	A	N1-C6-N6	9.28	124.17	118.60
1	AA	1649	A	N1-C6-N6	9.28	124.17	118.60
53	B5	2276	G	N1-C6-O6	9.28	125.47	119.90
53	B5	2747	A	N1-C6-N6	9.28	124.17	118.60
53	B5	593	A	N1-C6-N6	9.27	124.16	118.60
53	B5	3006	A	N1-C6-N6	9.27	124.17	118.60
51	B3	105	A	N1-C6-N6	9.27	124.16	118.60
53	B5	2948	C	O4'-C1'-N1	9.27	115.62	108.20
1	AA	1419	A	N1-C6-N6	9.27	124.16	118.60
53	B5	2291	A	N1-C6-N6	9.27	124.16	118.60
1	AA	1252	A	N1-C6-N6	9.27	124.16	118.60
53	B5	2222	A	N1-C6-N6	9.26	124.16	118.60
1	AA	892	A	N1-C6-N6	9.26	124.16	118.60
53	B5	1130	A	N1-C6-N6	9.26	124.16	118.60
1	AA	746	A	N1-C6-N6	9.26	124.16	118.60
53	B5	929	A	N1-C6-N6	9.26	124.16	118.60
1	AA	179	A	N1-C6-N6	9.26	124.16	118.60
1	AA	1747	A	N1-C6-N6	9.26	124.15	118.60
53	B5	1011	A	N1-C6-N6	9.26	124.15	118.60
53	B5	1065	A	N1-C6-N6	9.26	124.15	118.60
53	B5	2930	A	N1-C6-N6	9.25	124.15	118.60
53	B5	398	A	N1-C6-N6	9.25	124.15	118.60
53	B5	858	A	N1-C6-N6	9.25	124.15	118.60
52	B4	126	A	N1-C6-N6	9.24	124.15	118.60
53	B5	2940	A	N1-C6-N6	9.24	124.15	118.60
53	B5	2353	G	N1-C6-O6	9.24	125.44	119.90
1	AA	1357	A	N1-C6-N6	9.24	124.14	118.60
53	B5	913	A	N1-C6-N6	9.24	124.14	118.60
53	B5	2439	A	N1-C6-N6	9.24	124.14	118.60
53	B5	2841	G	N1-C6-O6	9.24	125.44	119.90
53	B5	2667	A	N1-C6-N6	9.23	124.14	118.60
1	AA	520	A	N1-C6-N6	9.23	124.14	118.60
53	B5	3314	A	N1-C6-N6	9.23	124.14	118.60
1	AA	1469	A	N1-C6-N6	9.23	124.14	118.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	1545	A	N1-C6-N6	9.23	124.14	118.60
53	B5	1521	G	N1-C6-O6	9.22	125.44	119.90
53	B5	1648	A	P-O3'-C3'	9.22	130.77	119.70
1	AA	1418	A	N1-C6-N6	9.22	124.13	118.60
1	AA	1029	A	N1-C6-N6	9.22	124.13	118.60
1	AA	1676	A	N1-C6-N6	9.22	124.13	118.60
53	B5	60	A	N1-C6-N6	9.22	124.13	118.60
53	B5	3295	A	N1-C6-N6	9.22	124.13	118.60
1	AA	1665	A	N1-C6-N6	9.22	124.13	118.60
53	B5	2657	A	N1-C6-N6	9.21	124.13	118.60
1	AA	1710	A	N1-C6-N6	9.21	124.13	118.60
53	B5	2929	C	O4'-C1'-N1	9.21	115.57	108.20
1	AA	601	A	N1-C6-N6	9.21	124.12	118.60
53	B5	1390	A	N1-C6-N6	9.21	124.12	118.60
53	B5	3122	A	N1-C6-N6	9.21	124.12	118.60
53	B5	647	A	N1-C6-N6	9.21	124.12	118.60
53	B5	709	A	N1-C6-N6	9.20	124.12	118.60
1	AA	331	A	N1-C6-N6	9.20	124.12	118.60
1	AA	844	A	N1-C6-N6	9.20	124.12	118.60
1	AA	1334	A	N1-C6-N6	9.20	124.12	118.60
1	AA	1795	A	N1-C6-N6	9.20	124.12	118.60
19	A7	52	U	N1-C2-N3	9.20	120.42	114.90
53	B5	99	A	N1-C6-N6	9.20	124.12	118.60
53	B5	1583	A	N1-C6-N6	9.20	124.12	118.60
53	B5	1799	A	N1-C6-N6	9.20	124.12	118.60
53	B5	395	A	N1-C6-N6	9.20	124.12	118.60
53	B5	1150	A	N1-C6-N6	9.20	124.12	118.60
53	B5	1263	A	N1-C6-N6	9.20	124.12	118.60
53	B5	2402	A	N1-C6-N6	9.20	124.12	118.60
51	B3	97	C	O4'-C1'-N1	9.20	115.56	108.20
53	B5	807	A	N1-C6-N6	9.20	124.12	118.60
53	B5	1240	A	N1-C6-N6	9.20	124.12	118.60
53	B5	1286	A	N1-C6-N6	9.19	124.11	118.60
53	B5	746	A	N1-C6-N6	9.19	124.11	118.60
53	B5	3307	A	N1-C6-N6	9.19	124.11	118.60
19	A7	71	G	N1-C6-O6	-9.19	114.39	119.90
51	B3	99	A	N1-C6-N6	9.19	124.11	118.60
53	B5	40	A	N1-C6-N6	9.19	124.11	118.60
53	B5	109	A	N1-C6-N6	9.19	124.11	118.60
53	B5	485	A	N1-C6-N6	9.19	124.11	118.60
1	AA	1189	A	N1-C6-N6	9.18	124.11	118.60
1	AA	1748	A	N1-C6-N6	9.18	124.11	118.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	1406	A	N1-C6-N6	9.18	124.11	118.60
1	AA	943	A	N1-C6-N6	9.18	124.11	118.60
53	B5	48	A	N1-C6-N6	9.18	124.11	118.60
53	B5	453	C	O4'-C1'-N1	9.18	115.54	108.20
1	AA	1084	A	N1-C6-N6	9.18	124.11	118.60
53	B5	1588	A	N1-C6-N6	9.18	124.11	118.60
53	B5	1913	A	N1-C6-N6	9.18	124.11	118.60
53	B5	1537	A	N1-C6-N6	9.17	124.10	118.60
1	AA	1192	A	N1-C6-N6	9.17	124.10	118.60
53	B5	1274	A	N1-C6-N6	9.17	124.10	118.60
53	B5	1346	G	N1-C6-O6	9.17	125.40	119.90
1	AA	881	A	N1-C6-N6	9.17	124.10	118.60
1	AA	1590	A	N1-C6-N6	9.17	124.10	118.60
19	A7	61	C	N3-C2-O2	-9.17	115.48	121.90
53	B5	199	A	N1-C6-N6	9.17	124.10	118.60
53	B5	2702	A	N1-C6-N6	9.17	124.10	118.60
53	B5	2233	A	N1-C6-N6	9.17	124.10	118.60
19	A7	69	U	O4'-C1'-N1	9.17	115.53	108.20
53	B5	1864	A	N1-C6-N6	9.17	124.10	118.60
1	AA	797	G	P-O5'-C5'	-9.16	106.24	120.90
53	B5	3034	C	O4'-C1'-N1	9.16	115.53	108.20
1	AA	545	A	N1-C6-N6	9.16	124.10	118.60
1	AA	566	C	O4'-C1'-N1	9.16	115.53	108.20
53	B5	175	C	O4'-C1'-N1	9.16	115.53	108.20
53	B5	2568	C	O4'-C1'-N1	9.16	115.53	108.20
1	AA	1293	A	N1-C6-N6	9.16	124.09	118.60
53	B5	130	A	N1-C6-N6	9.16	124.09	118.60
53	B5	1091	A	N1-C6-N6	9.15	124.09	118.60
53	B5	1301	A	N1-C6-N6	9.15	124.09	118.60
1	AA	762	A	N1-C6-N6	9.15	124.09	118.60
53	B5	1865	A	N1-C6-N6	9.15	124.09	118.60
1	AA	1574	A	N1-C6-N6	9.14	124.08	118.60
1	AA	1323	A	N1-C6-N6	9.14	124.08	118.60
53	B5	1384	U	O4'-C1'-N1	9.14	115.51	108.20
53	B5	2536	A	N1-C6-N6	9.14	124.08	118.60
53	B5	1612	A	N1-C6-N6	9.14	124.08	118.60
53	B5	1667	A	N1-C6-N6	9.14	124.08	118.60
53	B5	1369	A	N1-C6-N6	9.13	124.08	118.60
53	B5	1863	G	N1-C6-O6	9.13	125.38	119.90
53	B5	2680	A	N1-C6-N6	9.13	124.08	118.60
53	B5	711	A	P-O3'-C3'	9.13	130.66	119.70
53	B5	2911	A	N1-C6-N6	9.13	124.08	118.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	369	A	N1-C6-N6	9.13	124.08	118.60
53	B5	1859	A	N1-C6-N6	9.13	124.08	118.60
53	B5	3120	C	O4'-C1'-N1	9.13	115.50	108.20
1	AA	1557	A	N1-C6-N6	9.13	124.08	118.60
53	B5	2285	C	O4'-C1'-N1	9.13	115.50	108.20
53	B5	1771	C	O4'-C1'-N1	9.13	115.50	108.20
53	B5	2382	G	N1-C6-O6	9.13	125.38	119.90
53	B5	3005	A	N1-C6-N6	9.12	124.08	118.60
1	AA	1219	A	N1-C6-N6	9.12	124.07	118.60
53	B5	1053	A	N1-C6-N6	9.12	124.07	118.60
53	B5	1875	G	N1-C6-O6	9.12	125.37	119.90
1	AA	51	A	N1-C6-N6	9.12	124.07	118.60
53	B5	2372	A	N1-C6-N6	9.12	124.07	118.60
53	B5	3112	G	N1-C6-O6	9.12	125.37	119.90
51	B3	20	A	N1-C6-N6	9.12	124.07	118.60
53	B5	2213	A	N1-C6-N6	9.11	124.07	118.60
1	AA	809	A	N1-C6-N6	9.11	124.07	118.60
1	AA	1608	G	N1-C6-O6	9.11	125.37	119.90
53	B5	515	C	O4'-C1'-N1	9.11	115.49	108.20
53	B5	1477	A	N1-C6-N6	9.11	124.07	118.60
53	B5	2804	A	N1-C6-N6	9.11	124.07	118.60
53	B5	3103	A	N1-C6-N6	9.11	124.06	118.60
1	AA	1444	A	N1-C6-N6	9.11	124.06	118.60
53	B5	2494	A	N1-C6-N6	9.11	124.06	118.60
53	B5	63	A	N1-C6-N6	9.10	124.06	118.60
53	B5	967	A	N1-C6-N6	9.10	124.06	118.60
53	B5	3087	A	N1-C6-N6	9.10	124.06	118.60
1	AA	1004	A	N1-C6-N6	9.10	124.06	118.60
51	B3	64	A	N1-C6-N6	9.10	124.06	118.60
53	B5	1099	A	N1-C6-N6	9.10	124.06	118.60
1	AA	891	A	N1-C6-N6	9.09	124.05	118.60
53	B5	866	A	N1-C6-N6	9.09	124.05	118.60
53	B5	3342	A	N1-C6-N6	9.09	124.05	118.60
1	AA	464	A	N1-C6-N6	9.09	124.05	118.60
1	AA	1343	A	N1-C6-N6	9.09	124.05	118.60
1	AA	299	A	N1-C6-N6	9.08	124.05	118.60
53	B5	1047	A	N1-C6-N6	9.08	124.05	118.60
1	AA	481	A	N1-C6-N6	9.08	124.05	118.60
19	A7	29	A	O4'-C1'-N9	9.08	115.46	108.20
53	B5	3139	A	N1-C6-N6	9.08	124.05	118.60
1	AA	236	A	N1-C6-N6	9.08	124.05	118.60
1	AA	468	A	N1-C6-N6	9.08	124.05	118.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	35	A	N1-C6-N6	9.08	124.05	118.60
53	B5	2281	A	N1-C6-N6	9.08	124.05	118.60
19	A7	13	C	O4'-C1'-N1	9.08	115.46	108.20
51	B3	62	A	N1-C6-N6	9.08	124.05	118.60
1	AA	104	A	N1-C6-N6	9.07	124.04	118.60
53	B5	1433	A	N1-C6-N6	9.07	124.04	118.60
19	A7	71	G	C5-C6-N1	9.06	116.03	111.50
53	B5	962	A	N1-C6-N6	9.06	124.04	118.60
53	B5	1260	A	N1-C6-N6	9.06	124.04	118.60
53	B5	2182	A	N1-C6-N6	9.06	124.04	118.60
1	AA	400	A	P-O3'-C3'	9.06	130.57	119.70
1	AA	635	A	N1-C6-N6	9.06	124.04	118.60
53	B5	2244	A	N1-C6-N6	9.06	124.04	118.60
53	B5	3217	A	N1-C6-N6	9.06	124.04	118.60
53	B5	3130	A	N1-C6-N6	9.06	124.04	118.60
53	B5	2988	C	O4'-C1'-N1	9.06	115.44	108.20
53	B5	151	A	N1-C6-N6	9.05	124.03	118.60
53	B5	912	G	N1-C6-O6	9.05	125.33	119.90
1	AA	80	A	N1-C6-N6	9.04	124.03	118.60
1	AA	1705	A	N1-C6-N6	9.04	124.03	118.60
52	B4	104	A	N1-C6-N6	9.04	124.02	118.60
53	B5	622	A	N1-C6-N6	9.04	124.02	118.60
53	B5	1254	C	O4'-C1'-N1	9.04	115.43	108.20
53	B5	2853	A	N1-C6-N6	9.04	124.02	118.60
1	AA	580	A	N1-C6-N6	9.03	124.02	118.60
53	B5	993	A	N1-C6-N6	9.03	124.02	118.60
53	B5	951	A	N1-C6-N6	9.03	124.02	118.60
1	AA	256	A	N1-C6-N6	9.03	124.02	118.60
1	AA	518	A	N1-C6-N6	9.03	124.02	118.60
51	B3	13	A	N1-C6-N6	9.03	124.02	118.60
53	B5	1252	A	N1-C6-N6	9.03	124.02	118.60
1	AA	623	A	N1-C6-N6	9.03	124.02	118.60
1	AA	1179	A	N1-C6-N6	9.03	124.02	118.60
53	B5	266	A	N1-C6-N6	9.02	124.01	118.60
53	B5	351	A	N1-C6-N6	9.02	124.02	118.60
53	B5	977	U	P-O3'-C3'	9.02	130.53	119.70
53	B5	559	A	N1-C6-N6	9.02	124.01	118.60
53	B5	699	A	N1-C6-N6	9.02	124.01	118.60
53	B5	1437	C	O4'-C1'-N1	9.02	115.42	108.20
1	AA	1575	A	N1-C6-N6	9.02	124.01	118.60
19	A7	21	A	C4-C5-C6	-9.02	112.49	117.00
53	B5	205	C	O4'-C1'-N1	9.01	115.41	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1213	A	N1-C6-N6	9.01	124.01	118.60
53	B5	1797	A	N1-C6-N6	9.01	124.00	118.60
19	A7	45	G	C4-C5-C6	-9.01	113.40	118.80
53	B5	1685	C	O4'-C1'-N1	9.01	115.41	108.20
53	B5	2168	A	N1-C6-N6	9.01	124.00	118.60
1	AA	353	A	N1-C6-N6	9.00	124.00	118.60
1	AA	1204	A	N1-C6-N6	9.00	124.00	118.60
53	B5	3150	A	N1-C6-N6	9.00	124.00	118.60
1	AA	829	A	N1-C6-N6	9.00	124.00	118.60
1	AA	1322	A	N1-C6-N6	9.00	124.00	118.60
53	B5	1057	A	N1-C6-N6	9.00	124.00	118.60
53	B5	1930	A	N1-C6-N6	9.00	124.00	118.60
53	B5	1901	A	N1-C6-N6	9.00	124.00	118.60
53	B5	2142	A	N1-C6-N6	9.00	124.00	118.60
53	B5	2933	A	N1-C6-N6	9.00	124.00	118.60
1	AA	1372	A	N1-C6-N6	8.99	124.00	118.60
1	AA	1407	A	N1-C6-N6	8.99	124.00	118.60
1	AA	1479	C	O4'-C1'-N1	8.99	115.39	108.20
53	B5	674	G	N1-C6-O6	8.99	125.30	119.90
53	B5	1129	A	N1-C6-N6	8.99	124.00	118.60
53	B5	323	A	N1-C6-N6	8.99	124.00	118.60
53	B5	2461	A	N1-C6-N6	8.99	124.00	118.60
53	B5	2705	A	N1-C6-N6	8.99	124.00	118.60
53	B5	2546	C	O4'-C1'-N1	8.99	115.39	108.20
53	B5	1035	G	N1-C6-O6	8.99	125.29	119.90
1	AA	1066	A	N1-C6-N6	8.98	123.99	118.60
1	AA	1325	G	N1-C6-O6	8.98	125.29	119.90
53	B5	2305	G	N1-C6-O6	8.98	125.29	119.90
1	AA	1636	G	C4'-C3'-C2'	-8.98	93.62	102.60
19	A7	51	G	N3-C2-N2	-8.98	113.61	119.90
53	B5	1255	C	O4'-C1'-N1	8.98	115.38	108.20
53	B5	2356	A	N1-C6-N6	8.98	123.99	118.60
53	B5	2458	A	N1-C6-N6	8.98	123.99	118.60
53	B5	2709	C	O4'-C1'-N1	8.98	115.38	108.20
53	B5	848	A	N1-C6-N6	8.97	123.98	118.60
1	AA	1517	U	P-O3'-C3'	-8.97	108.93	119.70
1	AA	1595	A	N1-C6-N6	8.97	123.98	118.60
53	B5	2361	A	N1-C6-N6	8.97	123.98	118.60
51	B3	91	C	O4'-C1'-N1	8.97	115.38	108.20
53	B5	1449	A	N1-C6-N6	8.96	123.98	118.60
53	B5	1482	A	N1-C6-N6	8.96	123.98	118.60
53	B5	193	C	O4'-C1'-N1	8.96	115.37	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	438	A	N1-C6-N6	8.96	123.98	118.60
53	B5	1666	G	O4'-C1'-N9	8.96	115.37	108.20
53	B5	95	A	N1-C6-N6	8.96	123.98	118.60
1	AA	147	A	N1-C6-N6	8.96	123.97	118.60
53	B5	1874	A	N1-C6-N6	8.96	123.97	118.60
19	A7	13	C	N1-C2-O2	8.96	124.27	118.90
1	AA	990	G	N1-C6-O6	8.96	125.27	119.90
53	B5	1481	A	N1-C6-N6	8.96	123.97	118.60
53	B5	1839	A	N1-C6-N6	8.95	123.97	118.60
53	B5	88	A	N1-C6-N6	8.95	123.97	118.60
1	AA	103	A	N1-C6-N6	8.94	123.97	118.60
19	A7	36	A	C6-N1-C2	8.94	123.97	118.60
19	A7	52	U	N3-C2-O2	-8.95	115.94	122.20
53	B5	816	A	N1-C6-N6	8.95	123.97	118.60
53	B5	1525	G	N1-C6-O6	8.94	125.27	119.90
53	B5	1632	A	N1-C6-N6	8.94	123.97	118.60
53	B5	2107	A	N1-C6-N6	8.94	123.96	118.60
53	B5	366	A	N1-C6-N6	8.94	123.96	118.60
53	B5	1520	G	N1-C6-O6	8.93	125.26	119.90
51	B3	22	A	N1-C6-N6	8.93	123.96	118.60
53	B5	2196	C	O4'-C1'-N1	8.93	115.34	108.20
53	B5	3167	A	N1-C6-N6	8.93	123.96	118.60
1	AA	1521	G	O5'-P-OP2	8.93	121.41	110.70
53	B5	2696	A	N1-C6-N6	8.93	123.96	118.60
19	A7	69	U	C2-N3-C4	-8.93	121.64	127.00
53	B5	837	A	N1-C6-N6	8.93	123.96	118.60
53	B5	1317	A	N1-C6-N6	8.93	123.95	118.60
53	B5	801	A	N1-C6-N6	8.92	123.95	118.60
51	B3	115	A	N1-C6-N6	8.92	123.95	118.60
53	B5	209	A	N1-C6-N6	8.92	123.95	118.60
53	B5	1800	A	N1-C6-N6	8.92	123.95	118.60
53	B5	336	A	N1-C6-N6	8.91	123.95	118.60
1	AA	93	A	N1-C6-N6	8.91	123.95	118.60
53	B5	372	A	N1-C6-N6	8.91	123.95	118.60
53	B5	537	A	N1-C6-N6	8.91	123.95	118.60
53	B5	1899	G	N1-C6-O6	8.91	125.25	119.90
53	B5	3141	A	N1-C6-N6	8.91	123.95	118.60
51	B3	11	A	N1-C6-N6	8.91	123.94	118.60
53	B5	1546	A	N1-C6-N6	8.91	123.94	118.60
53	B5	2758	A	N1-C6-N6	8.91	123.94	118.60
53	B5	1363	A	N1-C6-N6	8.90	123.94	118.60
53	B5	3161	C	O4'-C1'-N1	8.90	115.32	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	306	A	N1-C6-N6	8.90	123.94	118.60
53	B5	839	C	O4'-C1'-N1	8.90	115.32	108.20
53	B5	3062	G	N1-C6-O6	8.90	125.24	119.90
52	B4	59	A	N1-C6-N6	8.90	123.94	118.60
53	B5	1275	C	O4'-C1'-N1	8.90	115.32	108.20
1	AA	316	A	N1-C6-N6	8.90	123.94	118.60
51	B3	46	A	N1-C6-N6	8.89	123.94	118.60
53	B5	3050	U	O4'-C1'-N1	8.89	115.31	108.20
19	A7	12	U	C5-C6-N1	-8.88	118.26	122.70
1	AA	535	A	N1-C6-N6	8.88	123.92	118.60
1	AA	1518	U	P-O3'-C3'	-8.87	109.06	119.70
53	B5	3053	G	N1-C6-O6	8.87	125.22	119.90
1	AA	592	A	N1-C6-N6	8.87	123.92	118.60
53	B5	728	G	N1-C6-O6	8.87	125.22	119.90
53	B5	2467	G	N1-C6-O6	8.87	125.22	119.90
1	AA	1764	A	N1-C6-N6	8.87	123.92	118.60
52	B4	65	A	N1-C6-N6	8.86	123.92	118.60
53	B5	2242	A	N1-C6-N6	8.86	123.92	118.60
53	B5	1366	A	N1-C6-N6	8.86	123.92	118.60
53	B5	1043	C	O4'-C1'-N1	8.86	115.28	108.20
53	B5	1500	G	N1-C6-O6	8.85	125.21	119.90
53	B5	1183	C	O4'-C1'-N1	8.85	115.28	108.20
53	B5	1776	G	N1-C6-O6	8.85	125.21	119.90
53	B5	156	G	N1-C6-O6	8.85	125.21	119.90
53	B5	1922	A	N1-C6-N6	8.85	123.91	118.60
53	B5	1334	U	O4'-C1'-N1	8.85	115.28	108.20
53	B5	1613	A	N1-C6-N6	8.85	123.91	118.60
1	AA	769	A	N1-C6-N6	8.84	123.91	118.60
19	A7	69	U	N3-C2-O2	-8.84	116.01	122.20
53	B5	108	A	N1-C6-N6	8.84	123.91	118.60
1	AA	436	A	N1-C6-N6	8.84	123.91	118.60
1	AA	1271	A	N1-C6-N6	8.84	123.91	118.60
19	A7	36	A	N1-C2-N3	-8.84	124.88	129.30
53	B5	2562	G	N1-C6-O6	8.84	125.20	119.90
52	B4	154	C	O4'-C1'-N1	8.83	115.27	108.20
53	B5	804	C	O4'-C1'-N1	8.83	115.27	108.20
53	B5	2449	A	N1-C6-N6	8.83	123.90	118.60
53	B5	3172	G	N1-C6-O6	8.83	125.20	119.90
1	AA	1127	G	N1-C6-O6	8.83	125.20	119.90
53	B5	1729	A	N1-C6-N6	8.83	123.90	118.60
53	B5	1779	C	O4'-C1'-N1	8.83	115.26	108.20
53	B5	2819	A	N1-C6-N6	8.83	123.90	118.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
19	A7	30	G	N9-C4-C5	8.82	108.93	105.40
53	B5	676	G	N1-C6-O6	8.82	125.19	119.90
53	B5	1826	C	O4'-C1'-N1	8.82	115.26	108.20
53	B5	1887	A	N1-C6-N6	8.82	123.89	118.60
53	B5	198	A	N1-C6-N6	8.82	123.89	118.60
53	B5	394	G	N1-C6-O6	8.82	125.19	119.90
53	B5	1217	A	N1-C6-N6	8.81	123.89	118.60
53	B5	1658	G	N1-C6-O6	8.81	125.18	119.90
53	B5	2780	A	N1-C6-N6	8.80	123.88	118.60
1	AA	1018	A	N1-C6-N6	8.80	123.88	118.60
53	B5	1184	A	N1-C6-N6	8.80	123.88	118.60
51	B3	104	C	O4'-C1'-N1	8.79	115.23	108.20
53	B5	66	A	N1-C6-N6	8.79	123.88	118.60
53	B5	729	C	O4'-C1'-N1	8.79	115.23	108.20
53	B5	1918	C	O4'-C1'-N1	8.78	115.23	108.20
1	AA	55	A	N1-C6-N6	8.78	123.87	118.60
53	B5	73	C	O4'-C1'-N1	8.78	115.22	108.20
1	AA	346	G	N1-C6-O6	8.78	125.17	119.90
1	AA	1198	A	N1-C6-N6	8.78	123.86	118.60
53	B5	2404	A	N1-C6-N6	8.78	123.86	118.60
1	AA	1335	C	O4'-C1'-N1	8.77	115.22	108.20
53	B5	20	A	N1-C6-N6	8.77	123.86	118.60
53	B5	127	G	N1-C6-O6	8.77	125.16	119.90
1	AA	460	A	N1-C6-N6	8.76	123.86	118.60
53	B5	290	G	N1-C6-O6	8.76	125.16	119.90
53	B5	2831	G	N1-C6-O6	8.76	125.16	119.90
1	AA	1520	U	C4'-C3'-C2'	8.76	111.36	102.60
19	A7	30	G	N3-C4-C5	-8.76	124.22	128.60
51	B3	40	C	O4'-C1'-N1	8.76	115.21	108.20
1	AA	1134	A	N1-C6-N6	8.75	123.85	118.60
1	AA	1019	A	N1-C6-N6	8.74	123.85	118.60
1	AA	1515	U	P-O3'-C3'	-8.74	109.21	119.70
53	B5	253	A	N1-C6-N6	8.74	123.85	118.60
53	B5	2525	G	N1-C6-O6	8.74	125.14	119.90
19	A7	18	G	O4'-C1'-N9	8.73	115.19	108.20
53	B5	1201	C	O4'-C1'-N1	8.73	115.19	108.20
1	AA	1627	G	N1-C6-O6	8.73	125.14	119.90
53	B5	2308	C	O4'-C1'-N1	8.73	115.19	108.20
53	B5	2686	A	N1-C6-N6	8.73	123.84	118.60
53	B5	2638	C	O4'-C1'-N1	8.73	115.18	108.20
1	AA	989	C	O4'-C1'-N1	8.73	115.18	108.20
53	B5	3261	C	O4'-C1'-N1	8.73	115.18	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	639	G	N1-C6-O6	8.72	125.13	119.90
53	B5	680	G	C5-C6-O6	-8.72	123.36	128.60
53	B5	3313	U	O4'-C1'-N1	8.72	115.18	108.20
53	B5	2459	A	N1-C6-N6	8.72	123.83	118.60
53	B5	2651	G	P-O3'-C3'	8.72	130.17	119.70
53	B5	15	C	O4'-C1'-N1	8.72	115.18	108.20
1	AA	872	G	N1-C6-O6	8.72	125.13	119.90
53	B5	321	C	O4'-C1'-N1	8.72	115.17	108.20
1	AA	315	A	N1-C6-N6	8.71	123.83	118.60
53	B5	159	A	N1-C6-N6	8.71	123.82	118.60
53	B5	1833	G	N1-C6-O6	8.71	125.12	119.90
53	B5	3038	U	O4'-C1'-N1	8.71	115.17	108.20
1	AA	451	A	N1-C6-N6	8.70	123.82	118.60
53	B5	1278	A	N1-C6-N6	8.70	123.82	118.60
1	AA	970	A	N1-C6-N6	8.70	123.82	118.60
53	B5	2150	G	N1-C6-O6	8.70	125.12	119.90
19	A7	36	A	N1-C6-N6	-8.69	113.39	118.60
53	B5	2856	G	N1-C6-O6	8.69	125.11	119.90
53	B5	2296	A	N1-C6-N6	8.69	123.81	118.60
53	B5	554	A	N1-C6-N6	8.69	123.81	118.60
53	B5	151	A	O4'-C1'-N9	8.68	115.14	108.20
53	B5	1828	A	N1-C6-N6	8.68	123.81	118.60
53	B5	2934	A	N1-C6-N6	8.68	123.81	118.60
1	AA	918	A	N1-C6-N6	8.68	123.81	118.60
1	AA	1197	G	N1-C6-O6	8.68	125.11	119.90
53	B5	384	A	N1-C6-N6	8.68	123.81	118.60
53	B5	440	A	N1-C6-N6	8.68	123.81	118.60
53	B5	1383	G	N1-C6-O6	8.68	125.11	119.90
53	B5	3088	G	N1-C6-O6	8.68	125.11	119.90
53	B5	2108	C	O4'-C1'-N1	8.68	115.14	108.20
53	B5	2313	A	N1-C6-N6	8.68	123.81	118.60
53	B5	2384	A	N1-C6-N6	8.67	123.80	118.60
53	B5	2976	A	N1-C6-N6	8.67	123.80	118.60
53	B5	3052	G	N1-C6-O6	8.67	125.10	119.90
17	AR	164	ASP	CB-CA-C	8.67	127.74	110.40
53	B5	2852	C	O4'-C1'-N1	8.67	115.13	108.20
1	AA	473	A	N1-C6-N6	8.66	123.80	118.60
53	B5	1947	G	O4'-C1'-N9	8.66	115.13	108.20
53	B5	2111	G	N1-C6-O6	8.66	125.10	119.90
1	AA	1110	A	N1-C6-N6	8.66	123.79	118.60
53	B5	1421	G	N1-C6-O6	8.66	125.09	119.90
53	B5	2708	C	O4'-C1'-N1	8.66	115.13	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	221	A	N1-C6-N6	8.65	123.79	118.60
53	B5	2309	A	N1-C6-N6	8.65	123.79	118.60
53	B5	3275	A	N1-C6-N6	8.65	123.79	118.60
1	AA	742	U	O4'-C1'-N1	8.65	115.12	108.20
1	AA	1024	A	N1-C6-N6	8.65	123.79	118.60
53	B5	1031	C	O4'-C1'-N1	8.65	115.12	108.20
53	B5	1318	A	N1-C6-N6	8.65	123.79	118.60
53	B5	2110	G	N1-C6-O6	8.65	125.09	119.90
1	AA	26	A	N1-C6-N6	8.64	123.79	118.60
1	AA	312	A	P-O3'-C3'	8.64	130.07	119.70
53	B5	259	C	O4'-C1'-N1	8.64	115.11	108.20
53	B5	1827	C	O4'-C1'-N1	8.64	115.11	108.20
53	B5	2420	C	O4'-C1'-N1	8.64	115.11	108.20
53	B5	2177	G	N1-C6-O6	8.63	125.08	119.90
51	B3	74	A	N1-C6-N6	8.62	123.78	118.60
53	B5	950	G	N1-C6-O6	8.62	125.07	119.90
19	A7	3	G	N3-C4-C5	-8.62	124.29	128.60
53	B5	2619	G	N1-C6-O6	8.62	125.07	119.90
53	B5	1257	C	O4'-C1'-N1	8.61	115.09	108.20
53	B5	595	A	N1-C6-N6	8.61	123.77	118.60
1	AA	1568	A	N1-C6-N6	8.61	123.77	118.60
53	B5	3152	U	O4'-C1'-N1	8.61	115.09	108.20
53	B5	1282	G	N1-C6-O6	8.61	125.06	119.90
53	B5	494	G	P-O3'-C3'	8.59	130.01	119.70
53	B5	2715	A	N1-C6-N6	8.59	123.75	118.60
53	B5	3225	C	O4'-C1'-N1	8.58	115.07	108.20
53	B5	1718	G	N1-C6-O6	8.58	125.05	119.90
53	B5	3061	G	N1-C6-O6	8.58	125.05	119.90
53	B5	132	C	O4'-C1'-N1	8.57	115.06	108.20
1	AA	47	A	P-O3'-C3'	8.57	129.99	119.70
1	AA	857	U	O4'-C1'-N1	8.57	115.06	108.20
53	B5	2126	A	N1-C6-N6	8.57	123.74	118.60
1	AA	865	A	P-O5'-C5'	-8.57	107.19	120.90
53	B5	815	G	N1-C6-O6	8.57	125.04	119.90
19	A7	44	A	C5'-C4'-C3'	-8.56	102.30	116.00
19	A7	57	G	C3'-C2'-C1'	8.56	108.35	101.50
53	B5	116	A	N1-C6-N6	8.56	123.74	118.60
53	B5	135	C	O4'-C1'-N1	8.56	115.05	108.20
53	B5	3004	C	O4'-C1'-N1	8.56	115.05	108.20
53	B5	2216	G	N1-C6-O6	8.55	125.03	119.90
53	B5	2192	C	O4'-C1'-N1	8.55	115.04	108.20
7	AH	61	ILE	CA-CB-CG1	-8.54	94.77	111.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1104	G	N1-C6-O6	8.54	125.03	119.90
52	B4	15	G	N1-C6-O6	8.54	125.03	119.90
53	B5	94	G	N1-C6-O6	8.54	125.02	119.90
51	B3	1	G	N1-C6-O6	8.54	125.02	119.90
53	B5	907	G	N1-C6-O6	8.54	125.02	119.90
53	B5	3037	U	O4'-C1'-N1	8.54	115.03	108.20
53	B5	1902	G	N1-C6-O6	8.53	125.02	119.90
53	B5	2117	A	N1-C6-N6	8.53	123.72	118.60
1	AA	1457	C	O4'-C1'-N1	8.53	115.02	108.20
53	B5	2265	C	O4'-C1'-N1	8.53	115.02	108.20
1	AA	474	A	N1-C6-N6	8.52	123.71	118.60
52	B4	112	U	O4'-C1'-N1	8.52	115.02	108.20
53	B5	2448	G	O4'-C1'-N9	8.52	115.02	108.20
53	B5	2823	G	N1-C6-O6	8.51	125.01	119.90
1	AA	992	A	N1-C6-N6	8.51	123.70	118.60
19	A7	47	U	C1'-O4'-C4'	-8.51	103.10	109.90
53	B5	1232	C	O4'-C1'-N1	8.51	115.00	108.20
53	B5	2906	C	O4'-C1'-N1	8.51	115.00	108.20
1	AA	309	C	O4'-C1'-N1	8.50	115.00	108.20
53	B5	477	A	N1-C6-N6	8.50	123.70	118.60
53	B5	1049	C	O4'-C1'-N1	8.50	115.00	108.20
53	B5	2237	C	O4'-C1'-N1	8.50	115.00	108.20
1	AA	973	A	N1-C6-N6	8.50	123.70	118.60
52	B4	11	C	O4'-C1'-N1	8.50	115.00	108.20
1	AA	1607	U	O4'-C1'-N1	8.49	115.00	108.20
1	AA	50	C	O4'-C1'-N1	8.49	114.99	108.20
1	AA	1499	C	O4'-C1'-N1	8.49	114.99	108.20
53	B5	3350	C	O4'-C1'-N1	8.49	114.99	108.20
1	AA	567	A	N1-C6-N6	8.48	123.69	118.60
53	B5	295	A	N1-C6-N6	8.48	123.69	118.60
53	B5	373	A	N1-C6-N6	8.48	123.69	118.60
53	B5	1140	G	N1-C6-O6	8.48	124.99	119.90
53	B5	2489	C	O4'-C1'-N1	8.48	114.98	108.20
53	B5	2331	C	O4'-C1'-N1	8.48	114.98	108.20
53	B5	2664	C	O4'-C1'-N1	8.48	114.98	108.20
53	B5	3116	G	N1-C6-O6	8.48	124.99	119.90
19	A7	62	A	C5-N7-C8	-8.47	99.66	103.90
53	B5	3125	U	O4'-C1'-N1	8.47	114.98	108.20
52	B4	142	C	O4'-C1'-N1	8.47	114.98	108.20
1	AA	168	A	N1-C6-N6	8.47	123.68	118.60
1	AA	867	G	C4'-C3'-C2'	8.47	111.07	102.60
53	B5	54	C	O4'-C1'-N1	8.47	114.97	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	1748	G	N1-C6-O6	8.47	124.98	119.90
53	B5	849	C	O4'-C1'-N1	8.47	114.97	108.20
53	B5	196	G	N1-C6-O6	8.46	124.98	119.90
53	B5	3129	A	N1-C6-N6	8.47	123.68	118.60
1	AA	568	G	N1-C6-O6	8.46	124.98	119.90
1	AA	1354	A	O4'-C1'-N9	8.46	114.97	108.20
19	A7	45	G	N7-C8-N9	8.46	117.33	113.10
1	AA	1548	A	P-O5'-C5'	-8.46	107.37	120.90
53	B5	637	C	O4'-C1'-N1	8.46	114.97	108.20
53	B5	2208	A	N1-C6-N6	8.46	123.67	118.60
53	B5	2476	C	O4'-C1'-N1	8.46	114.97	108.20
1	AA	1070	G	N1-C6-O6	8.45	124.97	119.90
1	AA	412	A	N1-C6-N6	8.45	123.67	118.60
19	A7	21	A	C5'-C4'-O4'	8.45	119.24	109.10
52	B4	141	C	O4'-C1'-N1	8.45	114.96	108.20
1	AA	1794	C	O4'-C1'-N1	8.45	114.96	108.20
53	B5	3274	G	N1-C6-O6	8.44	124.97	119.90
19	A7	24	G	C2-N3-C4	8.44	116.12	111.90
53	B5	2210	G	N1-C6-O6	8.43	124.96	119.90
53	B5	953	G	N1-C6-O6	8.43	124.95	119.90
52	B4	130	C	O4'-C1'-N1	8.42	114.94	108.20
19	A7	36	A	C8-N9-C4	-8.41	102.43	105.80
1	AA	856	A	N3-C4-C5	-8.41	120.91	126.80
52	B4	57	C	O4'-C1'-N1	8.41	114.93	108.20
53	B5	1117	G	N1-C6-O6	8.41	124.95	119.90
53	B5	2510	U	O4'-C1'-N1	8.41	114.93	108.20
53	B5	3112	G	C5-C6-O6	-8.41	123.55	128.60
1	AA	591	A	N1-C6-N6	8.41	123.64	118.60
19	A7	43	G	O4'-C1'-N9	8.41	114.92	108.20
53	B5	2299	A	N1-C6-N6	8.41	123.64	118.60
1	AA	587	C	O4'-C1'-N1	8.40	114.92	108.20
53	B5	1790	G	N1-C6-O6	8.40	124.94	119.90
53	B5	3091	A	N1-C6-N6	8.40	123.64	118.60
1	AA	176	C	O4'-C1'-N1	8.40	114.92	108.20
53	B5	1127	G	N1-C6-O6	8.40	124.94	119.90
1	AA	1071	G	N1-C6-O6	8.40	124.94	119.90
1	AA	1634	C	P-O3'-C3'	-8.40	109.62	119.70
53	B5	2263	C	O4'-C1'-N1	8.39	114.91	108.20
53	B5	2329	C	O4'-C1'-N1	8.39	114.91	108.20
53	B5	96	G	N1-C6-O6	8.39	124.93	119.90
53	B5	2928	C	O4'-C1'-N1	8.38	114.91	108.20
1	AA	1079	C	O4'-C1'-N1	8.38	114.90	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	64	G	N1-C6-O6	8.38	124.93	119.90
53	B5	2566	C	O4'-C1'-N1	8.38	114.90	108.20
53	B5	3338	C	O4'-C1'-N1	8.38	114.90	108.20
53	B5	1161	G	N1-C6-O6	8.38	124.93	119.90
53	B5	2630	C	O4'-C1'-N1	8.38	114.90	108.20
53	B5	3332	U	P-O3'-C3'	8.38	129.75	119.70
1	AA	319	U	O4'-C1'-N1	8.38	114.90	108.20
53	B5	133	U	P-O3'-C3'	8.38	129.75	119.70
53	B5	435	C	O4'-C1'-N1	8.38	114.90	108.20
1	AA	1425	G	N1-C6-O6	8.37	124.92	119.90
53	B5	695	C	O4'-C1'-N1	8.37	114.90	108.20
51	B3	88	G	N1-C6-O6	8.37	124.92	119.90
53	B5	2755	C	O4'-C1'-N1	8.37	114.89	108.20
53	B5	2853	A	O4'-C1'-N9	8.37	114.89	108.20
19	A7	71	G	C6-C5-N7	8.37	135.42	130.40
53	B5	826	G	N1-C6-O6	8.37	124.92	119.90
53	B5	2845	A	O4'-C1'-N9	8.37	114.89	108.20
53	B5	3033	A	N1-C6-N6	8.36	123.62	118.60
1	AA	1520	U	O4'-C4'-C3'	-8.36	95.64	104.00
1	AA	1093	C	O4'-C1'-N1	8.36	114.89	108.20
53	B5	678	G	N1-C6-O6	8.36	124.91	119.90
19	A7	27	C	C4'-C3'-C2'	-8.35	94.25	102.60
53	B5	2762	A	N1-C6-N6	8.35	123.61	118.60
53	B5	2332	A	N1-C6-N6	8.35	123.61	118.60
53	B5	1447	G	N1-C6-O6	8.35	124.91	119.90
1	AA	1587	C	O4'-C1'-N1	8.35	114.88	108.20
51	B3	75	G	C5-C6-O6	-8.34	123.59	128.60
53	B5	1917	C	O4'-C1'-N1	8.34	114.88	108.20
53	B5	1947	G	P-O3'-C3'	8.34	129.71	119.70
1	AA	625	C	O4'-C1'-N1	8.34	114.87	108.20
53	B5	1359	C	O4'-C1'-N1	8.34	114.87	108.20
53	B5	3248	C	O4'-C1'-N1	8.34	114.87	108.20
53	B5	787	G	N1-C6-O6	8.34	124.90	119.90
53	B5	1107	C	O4'-C1'-N1	8.34	114.87	108.20
53	B5	1735	G	N1-C6-O6	8.34	124.90	119.90
53	B5	1139	G	N1-C6-O6	8.34	124.90	119.90
53	B5	2845	A	N1-C6-N6	8.34	123.60	118.60
19	A7	43	G	N3-C2-N2	-8.33	114.07	119.90
53	B5	1487	G	N1-C6-O6	8.33	124.90	119.90
53	B5	1423	C	O4'-C1'-N1	8.33	114.87	108.20
53	B5	2812	C	O4'-C1'-N1	8.33	114.86	108.20
53	B5	3190	C	O4'-C1'-N1	8.33	114.86	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	68	C	O4'-C1'-N1	8.33	114.86	108.20
1	AA	866	G	P-O3'-C3'	-8.33	109.71	119.70
1	AA	1075	C	O4'-C1'-N1	8.32	114.86	108.20
53	B5	727	G	N1-C6-O6	8.32	124.89	119.90
53	B5	840	C	O4'-C1'-N1	8.32	114.86	108.20
53	B5	1414	G	N1-C6-O6	8.32	124.89	119.90
53	B5	2963	C	O4'-C1'-N1	8.32	114.85	108.20
1	AA	564	G	N1-C6-O6	8.31	124.89	119.90
19	A7	2	C	C5-C6-N1	-8.31	116.84	121.00
1	AA	1701	C	O4'-C1'-N1	8.31	114.85	108.20
53	B5	1393	A	N1-C6-N6	8.31	123.59	118.60
53	B5	2678	A	N1-C6-N6	8.31	123.59	118.60
1	AA	1521	G	N1-C2-N3	-8.31	118.91	123.90
52	B4	71	G	O4'-C1'-N9	8.31	114.85	108.20
53	B5	572	A	N1-C6-N6	8.31	123.58	118.60
53	B5	3102	G	N1-C6-O6	8.31	124.89	119.90
52	B4	137	C	O4'-C1'-N1	8.31	114.84	108.20
19	A7	59	U	C6-N1-C2	-8.30	116.02	121.00
53	B5	1420	C	O4'-C1'-N1	8.30	114.84	108.20
1	AA	858	G	N3-C4-N9	8.30	130.98	126.00
1	AA	1472	G	N1-C6-O6	8.30	124.88	119.90
53	B5	1387	G	N1-C6-O6	8.30	124.88	119.90
53	B5	609	G	N1-C6-O6	8.30	124.88	119.90
53	B5	1216	C	O4'-C1'-N1	8.30	114.84	108.20
53	B5	539	C	O4'-C1'-N1	8.30	114.84	108.20
53	B5	1653	G	N1-C6-O6	8.29	124.88	119.90
53	B5	1157	G	N1-C6-O6	8.29	124.88	119.90
53	B5	2651	G	N1-C6-O6	8.29	124.88	119.90
53	B5	2623	G	N1-C6-O6	8.29	124.87	119.90
53	B5	2245	C	O4'-C1'-N1	8.29	114.83	108.20
19	A7	44	A	N7-C8-N9	8.29	117.94	113.80
53	B5	2625	C	O4'-C1'-N1	8.29	114.83	108.20
53	B5	906	A	N1-C6-N6	8.29	123.57	118.60
53	B5	971	C	O4'-C1'-N1	8.28	114.83	108.20
1	AA	501	U	P-O3'-C3'	8.28	129.64	119.70
53	B5	809	G	N1-C6-O6	8.28	124.87	119.90
53	B5	1466	G	N1-C6-O6	8.28	124.87	119.90
53	B5	2496	C	O4'-C1'-N1	8.28	114.83	108.20
1	AA	1067	C	O4'-C1'-N1	8.28	114.82	108.20
1	AA	1277	G	N1-C6-O6	8.28	124.87	119.90
53	B5	2323	G	N1-C6-O6	8.28	124.87	119.90
53	B5	2682	C	O4'-C1'-N1	8.28	114.82	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
19	A7	15	G	C5-C6-O6	8.28	133.56	128.60
53	B5	1360	C	O4'-C1'-N1	8.27	114.82	108.20
53	B5	1365	G	N1-C6-O6	8.27	124.86	119.90
53	B5	2469	G	N1-C6-O6	8.27	124.86	119.90
51	B3	26	C	O4'-C1'-N1	8.27	114.81	108.20
51	B3	50	U	O4'-C1'-N1	8.27	114.81	108.20
52	B4	68	G	N1-C6-O6	8.27	124.86	119.90
51	B3	113	C	O4'-C1'-N1	8.26	114.81	108.20
53	B5	425	G	N1-C6-O6	8.26	124.86	119.90
53	B5	1900	A	P-O3'-C3'	8.26	129.61	119.70
53	B5	2151	C	O4'-C1'-N1	8.26	114.81	108.20
53	B5	2465	G	N1-C6-O6	8.26	124.86	119.90
53	B5	2276	G	C5-C6-O6	-8.26	123.65	128.60
53	B5	1190	A	N1-C6-N6	8.25	123.55	118.60
53	B5	2158	A	N1-C6-N6	8.25	123.55	118.60
53	B5	3228	C	O4'-C1'-N1	8.25	114.80	108.20
53	B5	3264	G	N1-C6-O6	8.25	124.85	119.90
1	AA	174	U	O4'-C1'-N1	8.25	114.80	108.20
53	B5	272	G	O4'-C1'-N9	8.25	114.80	108.20
19	A7	63	C	C1'-O4'-C4'	-8.24	103.31	109.90
53	B5	2297	U	O4'-C1'-N1	8.24	114.80	108.20
53	B5	1719	G	N1-C6-O6	8.24	124.84	119.90
53	B5	3030	G	N1-C6-O6	8.24	124.85	119.90
53	B5	824	C	O4'-C1'-N1	8.24	114.79	108.20
53	B5	1845	G	N1-C6-O6	8.24	124.84	119.90
53	B5	2545	C	O4'-C1'-N1	8.24	114.79	108.20
53	B5	3294	A	N1-C6-N6	8.24	123.54	118.60
53	B5	910	G	N1-C6-O6	8.24	124.84	119.90
53	B5	1186	G	N1-C6-O6	8.24	124.84	119.90
53	B5	1870	C	O4'-C1'-N1	8.24	114.79	108.20
53	B5	1251	A	O4'-C1'-N9	8.23	114.79	108.20
53	B5	968	G	N1-C6-O6	8.23	124.84	119.90
53	B5	696	C	O4'-C1'-N1	8.23	114.78	108.20
53	B5	522	A	N1-C6-N6	8.23	123.54	118.60
53	B5	2511	A	O4'-C1'-N9	8.23	114.78	108.20
19	A7	74	C	O4'-C1'-N1	8.22	114.78	108.20
53	B5	2704	A	N1-C6-N6	8.22	123.53	118.60
53	B5	2382	G	P-O3'-C3'	8.22	129.56	119.70
52	B4	78	G	N1-C6-O6	8.22	124.83	119.90
53	B5	2163	C	O4'-C1'-N1	8.22	114.78	108.20
1	AA	1581	A	N1-C6-N6	8.21	123.53	118.60
53	B5	714	G	N1-C6-O6	8.21	124.83	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	2783	U	O4'-C1'-N1	8.22	114.77	108.20
53	B5	2849	C	O4'-C1'-N1	8.21	114.77	108.20
1	AA	1638	C	OP1-P-OP2	-8.21	107.28	119.60
51	B3	48	U	O4'-C1'-N1	8.21	114.77	108.20
53	B5	2407	C	O4'-C1'-N1	8.21	114.77	108.20
1	AA	307	G	P-O3'-C3'	8.21	129.55	119.70
53	B5	312	C	O4'-C1'-N1	8.21	114.76	108.20
1	AA	976	A	N1-C6-N6	8.20	123.52	118.60
52	B4	150	G	N1-C6-O6	8.20	124.82	119.90
53	B5	258	G	N1-C6-O6	8.20	124.82	119.90
53	B5	1582	C	O4'-C1'-N1	8.20	114.76	108.20
53	B5	3367	C	O4'-C1'-N1	8.20	114.76	108.20
1	AA	1370	C	O4'-C1'-N1	8.19	114.75	108.20
53	B5	233	C	O4'-C1'-N1	8.19	114.75	108.20
53	B5	1119	C	O4'-C1'-N1	8.19	114.75	108.20
53	B5	1429	G	N1-C6-O6	8.19	124.81	119.90
53	B5	3101	G	N1-C6-O6	8.19	124.81	119.90
53	B5	3284	G	N1-C6-O6	8.19	124.81	119.90
53	B5	1219	C	O4'-C1'-N1	8.19	114.75	108.20
51	B3	18	C	O4'-C1'-N1	8.18	114.75	108.20
53	B5	484	C	O4'-C1'-N1	8.18	114.75	108.20
53	B5	1276	U	O4'-C1'-N1	8.18	114.75	108.20
1	AA	1637	C	P-O3'-C3'	-8.18	109.89	119.70
53	B5	1927	G	N1-C6-O6	8.18	124.81	119.90
53	B5	3098	G	N1-C6-O6	8.18	124.81	119.90
19	A7	43	G	N1-C2-N3	8.18	128.81	123.90
53	B5	928	C	O4'-C1'-N1	8.18	114.74	108.20
53	B5	1652	G	N1-C6-O6	8.18	124.81	119.90
1	AA	1201	C	O4'-C1'-N1	8.17	114.74	108.20
1	AA	1216	C	O4'-C1'-N1	8.17	114.74	108.20
53	B5	614	C	O4'-C1'-N1	8.17	114.74	108.20
1	AA	434	G	N1-C6-O6	8.17	124.80	119.90
1	AA	1064	C	O4'-C1'-N1	8.17	114.74	108.20
52	B4	129	C	O4'-C1'-N1	8.17	114.73	108.20
1	AA	969	A	N1-C6-N6	8.17	123.50	118.60
53	B5	1377	G	N1-C6-O6	8.17	124.80	119.90
1	AA	2	A	N1-C6-N6	8.16	123.50	118.60
1	AA	57	G	N1-C6-O6	8.16	124.80	119.90
53	B5	1508	C	O4'-C1'-N1	8.16	114.73	108.20
53	B5	2267	C	O4'-C1'-N1	8.16	114.73	108.20
53	B5	2526	C	O4'-C1'-N1	8.16	114.73	108.20
53	B5	2577	C	O4'-C1'-N1	8.16	114.73	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
51	B3	66	C	O4'-C1'-N1	8.16	114.73	108.20
1	AA	4	C	O4'-C1'-N1	8.16	114.72	108.20
53	B5	1243	G	N1-C6-O6	8.15	124.79	119.90
1	AA	1156	C	O4'-C1'-N1	8.15	114.72	108.20
1	AA	1400	C	O4'-C1'-N1	8.15	114.72	108.20
19	A7	31	A	O4'-C1'-N9	8.15	114.72	108.20
53	B5	1592	G	O4'-C1'-N9	8.15	114.72	108.20
53	B5	3290	G	N1-C6-O6	8.15	124.79	119.90
19	A7	3	G	N1-C6-O6	-8.15	115.01	119.90
53	B5	752	C	O4'-C1'-N1	8.15	114.72	108.20
53	B5	3084	C	O4'-C1'-N1	8.15	114.72	108.20
1	AA	796	A	OP1-P-OP2	-8.14	107.39	119.60
53	B5	2620	G	N1-C6-O6	8.14	124.79	119.90
53	B5	592	G	N1-C6-O6	8.14	124.78	119.90
53	B5	768	C	O4'-C1'-N1	8.14	114.71	108.20
53	B5	1464	G	N1-C6-O6	8.14	124.78	119.90
1	AA	298	C	O4'-C1'-N1	8.14	114.71	108.20
1	AA	1152	G	N1-C6-O6	8.13	124.78	119.90
53	B5	1285	G	N1-C6-O6	8.13	124.78	119.90
53	B5	1389	G	N1-C6-O6	8.13	124.78	119.90
53	B5	22	G	N1-C6-O6	8.13	124.78	119.90
53	B5	3128	G	C5-C6-O6	-8.13	123.72	128.60
1	AA	894	U	O4'-C1'-N1	8.13	114.70	108.20
19	A7	63	C	C6-N1-C2	-8.13	117.05	120.30
53	B5	574	U	O4'-C1'-N1	8.13	114.70	108.20
53	B5	2307	G	N1-C6-O6	8.13	124.78	119.90
53	B5	1340	G	N1-C6-O6	8.13	124.78	119.90
53	B5	2616	C	O4'-C1'-N1	8.12	114.70	108.20
1	AA	860	U	C4'-C3'-C2'	8.12	110.72	102.60
53	B5	1861	G	N1-C6-O6	8.12	124.77	119.90
53	B5	562	C	O4'-C1'-N1	8.12	114.70	108.20
53	B5	2583	C	O4'-C1'-N1	8.12	114.70	108.20
19	A7	71	G	N1-C2-N2	8.12	123.51	116.20
53	B5	460	C	O4'-C1'-N1	8.12	114.69	108.20
53	B5	1767	C	O4'-C1'-N1	8.12	114.69	108.20
53	B5	421	G	O4'-C1'-N9	8.11	114.69	108.20
53	B5	561	C	O4'-C1'-N1	8.11	114.69	108.20
53	B5	1068	C	O4'-C1'-N1	8.12	114.69	108.20
53	B5	2204	C	O4'-C1'-N1	8.12	114.69	108.20
53	B5	3208	C	O4'-C1'-N1	8.12	114.69	108.20
1	AA	1105	G	N1-C6-O6	8.11	124.77	119.90
19	A7	24	G	N3-C4-C5	-8.11	124.54	128.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	76	G	N1-C6-O6	8.11	124.77	119.90
1	AA	1241	G	N1-C6-O6	8.11	124.77	119.90
53	B5	407	A	N1-C6-N6	8.11	123.47	118.60
53	B5	1283	C	O4'-C1'-N1	8.11	114.69	108.20
19	A7	21	A	N9-C1'-C2'	-8.11	103.08	112.00
53	B5	274	G	N1-C6-O6	8.11	124.76	119.90
53	B5	1246	G	N1-C6-O6	8.11	124.76	119.90
53	B5	2927	C	O4'-C1'-N1	8.11	114.69	108.20
52	B4	25	G	N1-C6-O6	8.10	124.76	119.90
53	B5	240	U	O4'-C1'-N1	8.10	114.68	108.20
53	B5	1825	G	N1-C6-O6	8.10	124.76	119.90
53	B5	1403	C	O4'-C1'-N1	8.10	114.68	108.20
51	B3	8	G	N1-C6-O6	8.10	124.76	119.90
1	AA	1436	C	O4'-C1'-N1	8.09	114.67	108.20
53	B5	1794	G	N1-C6-O6	8.09	124.75	119.90
53	B5	155	G	N1-C6-O6	8.09	124.75	119.90
53	B5	1412	G	N1-C6-O6	8.09	124.75	119.90
53	B5	2618	G	N1-C6-O6	8.09	124.75	119.90
53	B5	19	U	O4'-C1'-N1	8.09	114.67	108.20
53	B5	1592	G	N1-C6-O6	8.09	124.75	119.90
53	B5	2202	C	O4'-C1'-N1	8.09	114.67	108.20
1	AA	1074	C	O4'-C1'-N1	8.08	114.66	108.20
53	B5	421	G	N1-C6-O6	8.08	124.75	119.90
53	B5	1069	C	O4'-C1'-N1	8.08	114.66	108.20
53	B5	1516	C	O4'-C1'-N1	8.08	114.67	108.20
1	AA	347	G	N1-C6-O6	8.08	124.75	119.90
1	AA	1371	C	O4'-C1'-N1	8.08	114.66	108.20
51	B3	15	C	O4'-C1'-N1	8.08	114.66	108.20
1	AA	172	C	O4'-C1'-N1	8.07	114.66	108.20
1	AA	339	C	O4'-C1'-N1	8.07	114.66	108.20
53	B5	3240	C	O4'-C1'-N1	8.07	114.66	108.20
53	B5	2945	G	N1-C6-O6	8.07	124.74	119.90
53	B5	2959	C	O4'-C1'-N1	8.07	114.66	108.20
1	AA	716	C	O4'-C1'-N1	8.07	114.65	108.20
53	B5	192	C	O4'-C1'-N1	8.07	114.65	108.20
1	AA	1020	C	O4'-C1'-N1	8.06	114.65	108.20
53	B5	111	C	O4'-C1'-N1	8.06	114.65	108.20
53	B5	427	C	O4'-C1'-N1	8.06	114.65	108.20
53	B5	1297	C	O4'-C1'-N1	8.06	114.65	108.20
53	B5	2830	G	O4'-C1'-N9	8.06	114.65	108.20
1	AA	310	C	O4'-C1'-N1	8.06	114.65	108.20
1	AA	1486	G	N1-C6-O6	8.06	124.74	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	332	C	O4'-C1'-N1	8.06	114.65	108.20
1	AA	184	C	O4'-C1'-N1	8.06	114.65	108.20
53	B5	237	G	N1-C6-O6	8.06	124.73	119.90
53	B5	1868	G	N1-C6-O6	8.06	124.73	119.90
53	B5	3312	U	O4'-C1'-N1	8.06	114.65	108.20
53	B5	1624	G	N1-C6-O6	8.06	124.73	119.90
53	B5	2483	G	N1-C6-O6	8.06	124.73	119.90
53	B5	2905	U	O4'-C1'-N1	8.06	114.65	108.20
51	B3	29	C	O4'-C1'-N1	8.06	114.64	108.20
1	AA	1171	C	O4'-C1'-N1	8.05	114.64	108.20
53	B5	1829	G	N1-C6-O6	8.05	124.73	119.90
1	AA	1401	C	O4'-C1'-N1	8.05	114.64	108.20
1	AA	645	C	O4'-C1'-N1	8.05	114.64	108.20
1	AA	1497	G	N1-C6-O6	8.05	124.73	119.90
53	B5	792	G	O4'-C1'-N9	8.05	114.64	108.20
53	B5	895	A	N1-C6-N6	8.05	123.43	118.60
53	B5	1928	G	N1-C6-O6	8.05	124.73	119.90
53	B5	3149	G	N1-C6-O6	8.05	124.73	119.90
1	AA	1336	C	O4'-C1'-N1	8.05	114.64	108.20
53	B5	1007	U	O4'-C1'-N1	8.05	114.64	108.20
53	B5	966	U	O4'-C1'-N1	8.05	114.64	108.20
53	B5	1846	C	O4'-C1'-N1	8.05	114.64	108.20
53	B5	2997	G	N1-C6-O6	8.05	124.73	119.90
1	AA	559	C	O4'-C1'-N1	8.04	114.63	108.20
53	B5	1376	C	O4'-C1'-N1	8.04	114.63	108.20
1	AA	644	C	O4'-C1'-N1	8.04	114.63	108.20
51	B3	95	C	O4'-C1'-N1	8.04	114.63	108.20
53	B5	2569	A	P-O3'-C3'	8.04	129.34	119.70
19	A7	9	A	C6-N1-C2	-8.04	113.78	118.60
53	B5	2303	A	O4'-C1'-N9	8.04	114.63	108.20
53	B5	2614	G	N1-C6-O6	8.04	124.72	119.90
53	B5	3289	G	N1-C6-O6	8.04	124.72	119.90
53	B5	1631	C	O4'-C1'-N1	8.03	114.63	108.20
53	B5	131	C	P-O3'-C3'	8.03	129.34	119.70
53	B5	404	G	N1-C6-O6	8.03	124.72	119.90
53	B5	1070	U	O4'-C1'-N1	8.03	114.62	108.20
53	B5	2961	G	N1-C6-O6	8.03	124.72	119.90
1	AA	627	C	O4'-C1'-N1	8.02	114.62	108.20
51	B3	39	C	O4'-C1'-N1	8.02	114.62	108.20
52	B4	105	A	N1-C6-N6	8.02	123.41	118.60
53	B5	1701	C	O4'-C1'-N1	8.02	114.62	108.20
53	B5	1725	C	O4'-C1'-N1	8.02	114.62	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	3092	C	O4'-C1'-N1	8.02	114.61	108.20
53	B5	2964	G	N1-C6-O6	8.02	124.71	119.90
53	B5	3241	G	N1-C6-O6	8.02	124.71	119.90
1	AA	773	C	O4'-C1'-N1	8.01	114.61	108.20
1	AA	777	C	O4'-C1'-N1	8.01	114.61	108.20
1	AA	1001	G	N1-C6-O6	8.01	124.71	119.90
51	B3	30	G	N1-C6-O6	8.01	124.71	119.90
53	B5	3193	C	O4'-C1'-N1	8.01	114.61	108.20
53	B5	2761	G	C5-C6-O6	-8.01	123.79	128.60
53	B5	2885	C	O4'-C1'-N1	8.01	114.61	108.20
19	A7	43	G	N3-C4-C5	-8.01	124.59	128.60
53	B5	139	G	N1-C6-O6	8.01	124.71	119.90
53	B5	3288	G	N1-C6-O6	8.01	124.70	119.90
53	B5	163	C	O4'-C1'-N1	8.01	114.61	108.20
53	B5	3022	G	N1-C6-O6	8.01	124.70	119.90
1	AA	674	C	O4'-C1'-N1	8.01	114.60	108.20
52	B4	144	G	N1-C6-O6	8.01	124.70	119.90
53	B5	260	C	O4'-C1'-N1	8.01	114.60	108.20
53	B5	281	G	N1-C6-O6	8.01	124.70	119.90
53	B5	1319	G	N1-C6-O6	8.01	124.70	119.90
53	B5	3197	G	N1-C6-O6	8.00	124.70	119.90
53	B5	463	C	O4'-C1'-N1	8.00	114.60	108.20
53	B5	1943	C	O4'-C1'-N1	8.00	114.60	108.20
1	AA	1416	G	N1-C6-O6	8.00	124.70	119.90
1	AA	784	C	O4'-C1'-N1	8.00	114.60	108.20
1	AA	961	C	O4'-C1'-N1	8.00	114.60	108.20
53	B5	1565	G	N1-C6-O6	8.00	124.70	119.90
1	AA	1274	G	N1-C6-O6	7.99	124.70	119.90
53	B5	1551	C	O4'-C1'-N1	7.99	114.59	108.20
53	B5	1654	A	O4'-C1'-N9	7.99	114.59	108.20
53	B5	3072	C	O4'-C1'-N1	7.99	114.59	108.20
53	B5	573	C	O4'-C1'-N1	7.99	114.59	108.20
53	B5	1628	C	O4'-C1'-N1	7.99	114.59	108.20
1	AA	708	C	O4'-C1'-N1	7.99	114.59	108.20
1	AA	1446	G	N1-C6-O6	7.99	124.69	119.90
53	B5	1764	U	O4'-C1'-N1	7.99	114.59	108.20
53	B5	105	C	O4'-C1'-N1	7.98	114.59	108.20
52	B4	110	C	O4'-C1'-N1	7.98	114.58	108.20
53	B5	190	U	O4'-C1'-N1	7.98	114.58	108.20
53	B5	1597	C	O4'-C1'-N1	7.98	114.58	108.20
53	B5	1947	G	N1-C6-O6	7.98	124.69	119.90
19	A7	8	U	C4-C5-C6	7.98	124.49	119.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	293	C	O4'-C1'-N1	7.98	114.58	108.20
1	AA	1373	C	O4'-C1'-N1	7.97	114.58	108.20
53	B5	849	C	C2-N1-C1'	7.97	127.57	118.80
53	B5	1668	G	N1-C6-O6	7.97	124.69	119.90
53	B5	2841	G	C5-C6-O6	-7.97	123.81	128.60
1	AA	1584	A	O4'-C1'-N9	7.97	114.58	108.20
52	B4	115	C	O4'-C1'-N1	7.97	114.58	108.20
53	B5	300	G	N1-C6-O6	7.97	124.68	119.90
53	B5	1733	G	N1-C6-O6	7.97	124.68	119.90
53	B5	2925	C	O4'-C1'-N1	7.97	114.58	108.20
53	B5	2710	C	O4'-C1'-N1	7.97	114.57	108.20
1	AA	1437	C	O4'-C1'-N1	7.96	114.57	108.20
1	AA	1520	U	C1'-C2'-O2'	7.96	134.49	110.60
53	B5	1499	C	O4'-C1'-N1	7.96	114.57	108.20
53	B5	2504	U	O4'-C1'-N1	7.96	114.57	108.20
53	B5	3369	G	N1-C6-O6	7.96	124.68	119.90
1	AA	703	G	N1-C6-O6	7.96	124.68	119.90
52	B4	106	C	O4'-C1'-N1	7.96	114.57	108.20
53	B5	1660	C	O4'-C1'-N1	7.96	114.57	108.20
53	B5	2189	U	O4'-C1'-N1	7.96	114.57	108.20
53	B5	1032	C	O4'-C1'-N1	7.96	114.57	108.20
53	B5	2277	C	O4'-C1'-N1	7.95	114.56	108.20
53	B5	1358	C	O4'-C1'-N1	7.95	114.56	108.20
53	B5	2490	C	O4'-C1'-N1	7.95	114.56	108.20
53	B5	881	C	O4'-C1'-N1	7.95	114.56	108.20
53	B5	2385	G	N1-C6-O6	7.95	124.67	119.90
53	B5	3060	C	O4'-C1'-N1	7.95	114.56	108.20
53	B5	3097	C	O4'-C1'-N1	7.95	114.56	108.20
1	AA	848	C	O4'-C1'-N1	7.95	114.56	108.20
53	B5	3366	G	N1-C6-O6	7.95	124.67	119.90
51	B3	93	U	O4'-C1'-N1	7.95	114.56	108.20
53	B5	2942	C	O4'-C1'-N1	7.95	114.56	108.20
53	B5	376	G	C5-C6-O6	-7.94	123.83	128.60
53	B5	767	U	O4'-C1'-N1	7.94	114.55	108.20
53	B5	1544	G	N1-C6-O6	7.94	124.67	119.90
53	B5	2836	C	O4'-C1'-N1	7.94	114.55	108.20
53	B5	148	G	N1-C6-O6	7.94	124.66	119.90
53	B5	1878	G	N1-C6-O6	7.94	124.66	119.90
53	B5	2622	C	O4'-C1'-N1	7.94	114.55	108.20
1	AA	115	G	N1-C6-O6	7.94	124.66	119.90
52	B4	133	G	O4'-C1'-N9	7.94	114.55	108.20
53	B5	225	C	O4'-C1'-N1	7.93	114.55	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	396	A	N1-C6-N6	7.93	123.36	118.60
53	B5	247	C	O4'-C1'-N1	7.93	114.55	108.20
53	B5	2444	C	O4'-C1'-N1	7.93	114.55	108.20
1	AA	867	G	C3'-C2'-C1'	-7.93	95.16	101.50
1	AA	717	C	O4'-C1'-N1	7.93	114.54	108.20
1	AA	491	C	O4'-C1'-N1	7.92	114.54	108.20
53	B5	1608	C	O4'-C1'-N1	7.92	114.54	108.20
52	B4	87	G	N1-C6-O6	7.92	124.65	119.90
53	B5	1518	U	O4'-C1'-N1	7.92	114.54	108.20
53	B5	1227	C	O4'-C1'-N1	7.92	114.53	108.20
53	B5	1357	G	N1-C6-O6	7.92	124.65	119.90
53	B5	3096	C	O4'-C1'-N1	7.92	114.53	108.20
1	AA	1770	C	O4'-C1'-N1	7.92	114.53	108.20
53	B5	1666	G	N1-C6-O6	7.92	124.65	119.90
53	B5	29	C	O4'-C1'-N1	7.91	114.53	108.20
53	B5	1067	U	O4'-C1'-N1	7.91	114.53	108.20
53	B5	2274	U	O4'-C1'-N1	7.91	114.53	108.20
53	B5	482	C	O4'-C1'-N1	7.91	114.53	108.20
53	B5	1416	C	O4'-C1'-N1	7.91	114.53	108.20
53	B5	2965	U	O4'-C1'-N1	7.91	114.53	108.20
53	B5	2218	G	N1-C6-O6	7.91	124.64	119.90
53	B5	2479	C	O4'-C1'-N1	7.91	114.53	108.20
53	B5	1131	G	N1-C6-O6	7.91	124.64	119.90
52	B4	151	C	O4'-C1'-N1	7.91	114.53	108.20
53	B5	3028	G	C5-C6-O6	-7.91	123.86	128.60
53	B5	927	C	O4'-C1'-N1	7.90	114.52	108.20
53	B5	1037	C	O4'-C1'-N1	7.90	114.52	108.20
53	B5	2350	C	O4'-C1'-N1	7.90	114.52	108.20
53	B5	2477	G	N1-C6-O6	7.90	124.64	119.90
53	B5	2542	U	O4'-C1'-N1	7.90	114.52	108.20
1	AA	442	C	O4'-C1'-N1	7.90	114.52	108.20
1	AA	974	C	O4'-C1'-N1	7.90	114.52	108.20
53	B5	383	G	N1-C6-O6	7.90	124.64	119.90
53	B5	1038	C	O4'-C1'-N1	7.90	114.52	108.20
53	B5	3311	C	O4'-C1'-N1	7.90	114.52	108.20
1	AA	1339	C	O4'-C1'-N1	7.90	114.52	108.20
53	B5	1905	G	N1-C6-O6	7.90	124.64	119.90
52	B4	51	G	N1-C6-O6	7.89	124.64	119.90
19	A7	13	C	C5'-C4'-O4'	7.89	118.57	109.10
53	B5	496	C	O4'-C1'-N1	7.89	114.51	108.20
53	B5	1563	C	O4'-C1'-N1	7.89	114.51	108.20
53	B5	1803	C	O4'-C1'-N1	7.89	114.52	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	2765	C	O4'-C1'-N1	7.89	114.51	108.20
1	AA	38	C	O4'-C1'-N1	7.89	114.51	108.20
1	AA	1528	C	O4'-C1'-N1	7.89	114.51	108.20
53	B5	206	G	N1-C6-O6	7.89	124.63	119.90
53	B5	1164	G	N1-C6-O6	7.89	124.63	119.90
53	B5	3233	C	O4'-C1'-N1	7.89	114.51	108.20
1	AA	1340	U	O4'-C1'-N1	7.89	114.51	108.20
53	B5	730	C	O4'-C1'-N1	7.89	114.51	108.20
1	AA	697	C	O4'-C1'-N1	7.88	114.51	108.20
53	B5	42	C	O4'-C1'-N1	7.88	114.51	108.20
53	B5	498	A	N1-C6-N6	7.88	123.33	118.60
53	B5	769	G	N1-C6-O6	7.88	124.63	119.90
53	B5	1658	G	C5-C6-O6	-7.88	123.87	128.60
53	B5	1497	C	O4'-C1'-N1	7.88	114.51	108.20
53	B5	2685	C	O4'-C1'-N1	7.88	114.50	108.20
1	AA	1675	C	O4'-C1'-N1	7.88	114.50	108.20
51	B3	71	C	O4'-C1'-N1	7.88	114.50	108.20
53	B5	346	C	O4'-C1'-N1	7.88	114.50	108.20
53	B5	3371	G	N1-C6-O6	7.88	124.63	119.90
1	AA	467	G	N1-C6-O6	7.88	124.63	119.90
53	B5	2707	C	O4'-C1'-N1	7.88	114.50	108.20
1	AA	658	C	O4'-C1'-N1	7.88	114.50	108.20
1	AA	1637	C	O5'-P-OP1	7.88	120.15	110.70
1	AA	735	C	O4'-C1'-N1	7.87	114.50	108.20
53	B5	1838	G	N1-C6-O6	7.87	124.62	119.90
53	B5	2826	U	P-O3'-C3'	7.87	129.15	119.70
53	B5	795	G	N1-C6-O6	7.87	124.62	119.90
53	B5	2646	C	O4'-C1'-N1	7.87	114.50	108.20
1	AA	1080	G	N1-C6-O6	7.87	124.62	119.90
53	B5	2862	U	O4'-C1'-N1	7.87	114.50	108.20
53	B5	1422	G	O4'-C1'-N9	7.87	114.50	108.20
52	B4	45	C	O4'-C1'-N1	7.87	114.49	108.20
1	AA	393	C	O4'-C1'-N1	7.86	114.49	108.20
19	A7	62	A	C4-C5-C6	-7.86	113.07	117.00
51	B3	57	C	O4'-C1'-N1	7.86	114.49	108.20
19	A7	75	C	N1-C2-N3	7.86	124.70	119.20
37	BM	148	LYS	C-N-CA	-7.86	102.05	121.70
53	B5	2893	C	O4'-C1'-N1	7.86	114.49	108.20
53	B5	158	G	N1-C6-O6	7.86	124.62	119.90
53	B5	480	C	O4'-C1'-N1	7.86	114.49	108.20
53	B5	2966	G	N1-C6-O6	7.86	124.62	119.90
1	AA	1616	C	O4'-C1'-N1	7.86	114.48	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1707	C	O4'-C1'-N1	7.86	114.48	108.20
53	B5	1024	G	N1-C6-O6	7.86	124.61	119.90
53	B5	1209	G	N1-C6-O6	7.86	124.61	119.90
53	B5	2355	G	N1-C6-O6	7.86	124.61	119.90
53	B5	2913	C	O4'-C1'-N1	7.86	114.49	108.20
53	B5	47	C	O4'-C1'-N1	7.86	114.48	108.20
53	B5	242	C	O4'-C1'-N1	7.86	114.48	108.20
53	B5	2221	G	N1-C6-O6	7.86	124.61	119.90
1	AA	1453	G	N1-C6-O6	7.85	124.61	119.90
53	B5	2562	G	C5-C6-O6	-7.85	123.89	128.60
53	B5	2666	C	O4'-C1'-N1	7.85	114.48	108.20
53	B5	2389	C	O4'-C1'-N1	7.85	114.48	108.20
1	AA	500	C	O4'-C1'-N1	7.85	114.48	108.20
1	AA	1050	G	O4'-C1'-N9	7.85	114.48	108.20
53	B5	359	U	O4'-C1'-N1	7.85	114.48	108.20
53	B5	670	C	O4'-C1'-N1	7.85	114.48	108.20
53	B5	1476	G	N1-C6-O6	7.85	124.61	119.90
53	B5	2422	C	O4'-C1'-N1	7.85	114.48	108.20
53	B5	3242	G	N1-C6-O6	7.85	124.61	119.90
53	B5	1756	C	O4'-C1'-N1	7.85	114.48	108.20
53	B5	1450	G	N1-C6-O6	7.84	124.61	119.90
53	B5	3333	G	N1-C6-O6	7.84	124.61	119.90
53	B5	489	C	O4'-C1'-N1	7.84	114.47	108.20
19	A7	28	C	C4-C5-C6	-7.84	113.48	117.40
53	B5	1042	U	O4'-C1'-N1	7.84	114.47	108.20
53	B5	1379	G	N1-C6-O6	7.84	124.61	119.90
53	B5	2512	C	O4'-C1'-N1	7.84	114.47	108.20
53	B5	2567	C	O4'-C1'-N1	7.84	114.47	108.20
1	AA	1584	A	O4'-C4'-C3'	-7.84	96.16	104.00
53	B5	1292	C	O4'-C1'-N1	7.84	114.47	108.20
53	B5	120	G	N1-C6-O6	7.84	124.60	119.90
53	B5	457	C	O4'-C1'-N1	7.84	114.47	108.20
53	B5	1158	A	N1-C6-N6	7.83	123.30	118.60
52	B4	103	G	N1-C6-O6	7.83	124.60	119.90
53	B5	753	C	O4'-C1'-N1	7.83	114.47	108.20
53	B5	861	C	O4'-C1'-N1	7.83	114.47	108.20
53	B5	1244	A	O4'-C1'-N9	7.83	114.47	108.20
53	B5	1549	U	O4'-C1'-N1	7.83	114.47	108.20
1	AA	230	C	O4'-C1'-N1	7.83	114.47	108.20
1	AA	465	G	N1-C6-O6	7.83	124.60	119.90
1	AA	1550	U	C3'-C2'-C1'	-7.83	95.23	101.50
53	B5	1185	C	O4'-C1'-N1	7.83	114.47	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	1441	G	N1-C6-O6	7.83	124.60	119.90
1	AA	286	C	O4'-C1'-N1	7.83	114.46	108.20
53	B5	92	G	N1-C6-O6	7.83	124.60	119.90
53	B5	948	C	O4'-C1'-N1	7.83	114.46	108.20
53	B5	2156	C	O4'-C1'-N1	7.83	114.46	108.20
1	AA	1043	G	N1-C6-O6	7.83	124.60	119.90
53	B5	597	G	N1-C6-O6	7.83	124.59	119.90
53	B5	2548	C	O4'-C1'-N1	7.83	114.46	108.20
53	B5	1422	G	N1-C6-O6	7.82	124.59	119.90
53	B5	1222	G	N1-C6-O6	7.82	124.59	119.90
1	AA	883	C	O4'-C1'-N1	7.82	114.46	108.20
19	A7	7	U	N3-C2-O2	-7.82	116.73	122.20
52	B4	47	C	O4'-C1'-N1	7.82	114.45	108.20
19	A7	20	G	N1-C6-O6	-7.82	115.21	119.90
53	B5	218	G	N1-C6-O6	7.82	124.59	119.90
53	B5	667	C	O4'-C1'-N1	7.82	114.45	108.20
53	B5	2828	G	N1-C6-O6	7.82	124.59	119.90
51	B3	110	C	O4'-C1'-N1	7.82	114.45	108.20
53	B5	499	G	N1-C6-O6	7.81	124.59	119.90
53	B5	2286	U	O4'-C1'-N1	7.81	114.45	108.20
1	AA	637	C	O4'-C1'-N1	7.81	114.45	108.20
53	B5	788	C	O4'-C1'-N1	7.81	114.45	108.20
53	B5	1869	C	O4'-C1'-N1	7.81	114.45	108.20
19	A7	22	G	C5'-C4'-O4'	7.81	118.47	109.10
53	B5	1665	C	O4'-C1'-N1	7.81	114.45	108.20
19	A7	48	C	O4'-C1'-N1	7.81	114.45	108.20
53	B5	613	G	N1-C6-O6	7.81	124.58	119.90
1	AA	1409	G	N1-C6-O6	7.80	124.58	119.90
52	B4	24	G	N1-C6-O6	7.80	124.58	119.90
53	B5	765	C	O4'-C1'-N1	7.80	114.44	108.20
1	AA	1009	C	O4'-C1'-N1	7.80	114.44	108.20
19	A7	66	A	N9-C4-C5	7.80	108.92	105.80
53	B5	1076	C	O4'-C1'-N1	7.80	114.44	108.20
53	B5	2579	G	N1-C6-O6	7.80	124.58	119.90
53	B5	2916	U	O4'-C1'-N1	7.80	114.44	108.20
1	AA	372	G	N1-C6-O6	7.80	124.58	119.90
1	AA	1518	U	O4'-C4'-C3'	-7.80	96.20	104.00
53	B5	1578	C	O4'-C1'-N1	7.80	114.44	108.20
53	B5	2810	C	O4'-C1'-N1	7.80	114.44	108.20
1	AA	1352	C	O4'-C1'-N1	7.80	114.44	108.20
19	A7	7	U	C5-C6-N1	-7.80	118.80	122.70
53	B5	957	C	O4'-C1'-N1	7.80	114.44	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	3132	C	O4'-C1'-N1	7.80	114.44	108.20
53	B5	3162	C	O4'-C1'-N1	7.80	114.44	108.20
53	B5	818	C	O4'-C1'-N1	7.79	114.44	108.20
53	B5	2544	U	O4'-C1'-N1	7.79	114.44	108.20
53	B5	908	G	N1-C6-O6	7.79	124.58	119.90
53	B5	3213	G	N1-C6-O6	7.79	124.58	119.90
53	B5	1333	C	O4'-C1'-N1	7.79	114.43	108.20
53	B5	1395	G	N1-C6-O6	7.79	124.57	119.90
53	B5	1586	G	N1-C6-O6	7.79	124.57	119.90
53	B5	3364	C	O4'-C1'-N1	7.79	114.43	108.20
1	AA	408	C	O4'-C1'-N1	7.79	114.43	108.20
1	AA	490	C	O4'-C1'-N1	7.79	114.43	108.20
53	B5	959	C	O4'-C1'-N1	7.79	114.43	108.20
53	B5	1397	C	O4'-C1'-N1	7.79	114.43	108.20
53	B5	1492	G	N1-C6-O6	7.79	124.57	119.90
53	B5	2879	C	O4'-C1'-N1	7.79	114.43	108.20
53	B5	2337	C	O4'-C1'-N1	7.79	114.43	108.20
1	AA	1036	C	O4'-C1'-N1	7.79	114.43	108.20
53	B5	945	C	O4'-C1'-N1	7.79	114.43	108.20
53	B5	3309	G	N1-C6-O6	7.79	124.57	119.90
52	B4	39	G	N1-C6-O6	7.78	124.57	119.90
53	B5	142	C	O4'-C1'-N1	7.78	114.43	108.20
1	AA	431	C	O4'-C1'-N1	7.78	114.42	108.20
53	B5	1116	G	N1-C6-O6	7.78	124.57	119.90
1	AA	189	C	O4'-C1'-N1	7.78	114.42	108.20
53	B5	785	G	N1-C6-O6	7.78	124.57	119.90
53	B5	2744	U	O4'-C1'-N1	7.78	114.42	108.20
53	B5	1770	G	N1-C6-O6	7.78	124.57	119.90
53	B5	271	C	O4'-C1'-N1	7.78	114.42	108.20
1	AA	613	G	N1-C6-O6	7.78	124.56	119.90
1	AA	926	C	O4'-C1'-N1	7.78	114.42	108.20
52	B4	95	G	O4'-C1'-N9	7.78	114.42	108.20
53	B5	455	C	O4'-C1'-N1	7.78	114.42	108.20
53	B5	638	C	O4'-C1'-N1	7.77	114.42	108.20
53	B5	1316	C	O4'-C1'-N1	7.77	114.42	108.20
53	B5	3285	C	O4'-C1'-N1	7.77	114.42	108.20
1	AA	1578	C	O4'-C1'-N1	7.77	114.42	108.20
53	B5	69	C	O4'-C1'-N1	7.77	114.42	108.20
53	B5	576	C	O4'-C1'-N1	7.77	114.42	108.20
53	B5	582	G	O4'-C1'-N9	7.77	114.42	108.20
53	B5	2370	G	N1-C6-O6	7.77	124.56	119.90
53	B5	3390	G	N1-C6-O6	7.77	124.56	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	1364	C	O4'-C1'-N1	7.77	114.42	108.20
1	AA	405	C	O4'-C1'-N1	7.77	114.41	108.20
1	AA	726	C	O4'-C1'-N1	7.77	114.41	108.20
53	B5	82	C	O4'-C1'-N1	7.77	114.41	108.20
53	B5	2478	C	O4'-C1'-N1	7.77	114.41	108.20
53	B5	3053	G	C5-C6-O6	-7.77	123.94	128.60
1	AA	1749	C	O4'-C1'-N1	7.76	114.41	108.20
19	A7	67	A	O4'-C1'-N9	7.76	114.41	108.20
53	B5	1693	C	O4'-C1'-N1	7.76	114.41	108.20
19	A7	42	G	N3-C4-C5	-7.76	124.72	128.60
53	B5	3207	C	O4'-C1'-N1	7.76	114.41	108.20
1	AA	1200	A	N3-C4-C5	-7.76	121.37	126.80
52	B4	8	C	O4'-C1'-N1	7.76	114.41	108.20
1	AA	1704	C	O4'-C1'-N1	7.76	114.41	108.20
52	B4	117	C	O4'-C1'-N1	7.76	114.41	108.20
53	B5	81	C	O4'-C1'-N1	7.76	114.41	108.20
53	B5	3235	C	O4'-C1'-N1	7.76	114.41	108.20
53	B5	1562	C	O4'-C1'-N1	7.75	114.40	108.20
53	B5	3063	C	O4'-C1'-N1	7.75	114.40	108.20
1	AA	955	C	O4'-C1'-N1	7.75	114.40	108.20
53	B5	525	C	O4'-C1'-N1	7.75	114.40	108.20
53	B5	757	C	O4'-C1'-N1	7.75	114.40	108.20
53	B5	921	A	N1-C6-N6	7.75	123.25	118.60
53	B5	1426	C	O4'-C1'-N1	7.75	114.40	108.20
1	AA	696	C	O4'-C1'-N1	7.75	114.40	108.20
53	B5	1132	C	O4'-C1'-N1	7.75	114.40	108.20
53	B5	1734	G	N1-C6-O6	7.75	124.55	119.90
53	B5	3257	C	O4'-C1'-N1	7.75	114.40	108.20
1	AA	1608	G	C5-C6-O6	-7.75	123.95	128.60
53	B5	2351	U	O4'-C1'-N1	7.75	114.40	108.20
1	AA	190	C	O4'-C1'-N1	7.75	114.40	108.20
1	AA	1563	C	O4'-C1'-N1	7.75	114.40	108.20
53	B5	886	C	O4'-C1'-N1	7.75	114.40	108.20
1	AA	530	C	O4'-C1'-N1	7.75	114.40	108.20
1	AA	898	G	N1-C6-O6	7.75	124.55	119.90
53	B5	880	G	N1-C6-O6	7.75	124.55	119.90
53	B5	3340	G	N1-C6-O6	7.75	124.55	119.90
19	A7	19	G	C4-C5-N7	7.74	113.90	110.80
53	B5	2475	G	O4'-C1'-N9	7.74	114.39	108.20
53	B5	2624	G	N1-C6-O6	7.74	124.55	119.90
1	AA	827	C	O4'-C1'-N1	7.74	114.39	108.20
1	AA	489	C	O4'-C1'-N1	7.74	114.39	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
19	A7	13	C	C3'-C2'-C1'	7.74	107.69	101.50
19	A7	19	G	C8-N9-C4	-7.74	103.31	106.40
53	B5	3015	G	N1-C6-O6	7.74	124.54	119.90
53	B5	74	G	N1-C6-O6	7.73	124.54	119.90
53	B5	1162	U	O4'-C1'-N1	7.73	114.39	108.20
53	B5	2452	G	N1-C6-O6	7.73	124.54	119.90
1	AA	688	G	N1-C6-O6	7.73	124.54	119.90
1	AA	1762	C	O4'-C1'-N1	7.73	114.38	108.20
53	B5	586	C	O4'-C1'-N1	7.73	114.39	108.20
53	B5	1816	A	O4'-C1'-N9	7.73	114.38	108.20
53	B5	2600	C	O4'-C1'-N1	7.73	114.39	108.20
1	AA	54	C	O4'-C1'-N1	7.73	114.38	108.20
53	B5	2582	C	O4'-C1'-N1	7.73	114.38	108.20
1	AA	191	C	O4'-C1'-N1	7.73	114.38	108.20
1	AA	1362	C	O4'-C1'-N1	7.73	114.38	108.20
19	A7	13	C	N3-C2-O2	-7.73	116.49	121.90
53	B5	3337	G	N1-C6-O6	7.73	124.54	119.90
1	AA	1768	U	O4'-C1'-N1	7.73	114.38	108.20
1	AA	1450	U	O4'-C1'-N1	7.72	114.38	108.20
1	AA	1461	C	O4'-C1'-N1	7.72	114.38	108.20
1	AA	1638	C	P-O5'-C5'	-7.72	108.54	120.90
53	B5	2899	C	O4'-C1'-N1	7.72	114.38	108.20
1	AA	1511	G	N1-C6-O6	7.72	124.53	119.90
52	B4	76	C	O4'-C1'-N1	7.72	114.38	108.20
53	B5	1303	A	O4'-C1'-N9	7.72	114.38	108.20
53	B5	1520	G	C5-C6-O6	-7.72	123.97	128.60
1	AA	1275	C	O4'-C1'-N1	7.72	114.38	108.20
53	B5	1747	G	N1-C6-O6	7.72	124.53	119.90
53	B5	2594	C	O4'-C1'-N1	7.72	114.38	108.20
1	AA	354	C	O4'-C1'-N1	7.72	114.38	108.20
1	AA	1125	C	O4'-C1'-N1	7.72	114.37	108.20
1	AA	1790	G	N1-C6-O6	7.72	124.53	119.90
52	B4	21	C	O4'-C1'-N1	7.72	114.37	108.20
1	AA	402	C	O4'-C1'-N1	7.71	114.37	108.20
1	AA	342	C	O4'-C1'-N1	7.71	114.37	108.20
53	B5	1573	G	N1-C6-O6	7.71	124.53	119.90
53	B5	1781	C	O4'-C1'-N1	7.71	114.37	108.20
1	AA	1651	C	O4'-C1'-N1	7.71	114.37	108.20
53	B5	963	G	N1-C6-O6	7.71	124.53	119.90
1	AA	765	G	O4'-C1'-N9	7.71	114.37	108.20
1	AA	1642	C	O4'-C1'-N1	7.71	114.37	108.20
53	B5	415	G	N1-C6-O6	7.71	124.53	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1613	C	O4'-C1'-N1	7.71	114.37	108.20
53	B5	2518	C	O4'-C1'-N1	7.71	114.37	108.20
53	B5	2835	U	O4'-C1'-N1	7.71	114.36	108.20
53	B5	3315	G	N1-C6-O6	7.71	124.52	119.90
53	B5	3355	U	O4'-C1'-N1	7.71	114.36	108.20
53	B5	291	C	O4'-C1'-N1	7.71	114.36	108.20
53	B5	2876	C	O4'-C1'-N1	7.71	114.36	108.20
53	B5	3148	U	O4'-C1'-N1	7.71	114.36	108.20
53	B5	379	C	O4'-C1'-N1	7.70	114.36	108.20
53	B5	1675	G	N1-C6-O6	7.70	124.52	119.90
53	B5	2729	U	O4'-C1'-N1	7.70	114.36	108.20
53	B5	2793	G	N1-C6-O6	7.70	124.52	119.90
46	BV	47	TYR	CB-CG-CD2	7.70	125.62	121.00
1	AA	501	U	N1-C2-N3	7.70	119.52	114.90
1	AA	1253	U	O4'-C1'-N1	7.70	114.36	108.20
19	A7	66	A	C4-C5-N7	-7.70	106.85	110.70
53	B5	644	G	N1-C6-O6	7.70	124.52	119.90
1	AA	1169	G	N1-C6-O6	7.69	124.52	119.90
1	AA	425	A	O4'-C1'-N9	7.69	114.36	108.20
53	B5	1472	U	O4'-C1'-N1	7.69	114.35	108.20
53	B5	1793	C	O4'-C1'-N1	7.69	114.35	108.20
1	AA	30	G	N1-C6-O6	7.69	124.51	119.90
1	AA	53	G	N1-C6-O6	7.69	124.51	119.90
53	B5	1596	C	O4'-C1'-N1	7.69	114.35	108.20
1	AA	730	G	N1-C6-O6	7.69	124.51	119.90
53	B5	1392	G	N1-C6-O6	7.69	124.51	119.90
53	B5	1852	G	N1-C6-O6	7.69	124.51	119.90
1	AA	1296	G	N1-C6-O6	7.69	124.51	119.90
1	AA	398	G	N1-C6-O6	7.68	124.51	119.90
1	AA	1775	G	N1-C6-O6	7.68	124.51	119.90
53	B5	140	C	O4'-C1'-N1	7.68	114.35	108.20
1	AA	986	G	N1-C6-O6	7.68	124.51	119.90
1	AA	1493	C	O4'-C1'-N1	7.68	114.34	108.20
53	B5	3221	C	O4'-C1'-N1	7.68	114.34	108.20
53	B5	3386	G	N1-C6-O6	7.68	124.51	119.90
1	AA	1098	G	N1-C6-O6	7.68	124.51	119.90
1	AA	1703	C	O4'-C1'-N1	7.68	114.34	108.20
53	B5	1277	C	O4'-C1'-N1	7.68	114.34	108.20
1	AA	1776	G	N1-C6-O6	7.68	124.51	119.90
53	B5	1655	G	N1-C6-O6	7.68	124.51	119.90
19	A7	18	G	N3-C4-N9	-7.67	121.39	126.00
53	B5	392	G	N1-C6-O6	7.67	124.50	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	3135	U	O4'-C1'-N1	7.67	114.34	108.20
53	B5	3360	C	O4'-C1'-N1	7.67	114.34	108.20
53	B5	633	C	O4'-C1'-N1	7.67	114.34	108.20
19	A7	5	A	C4'-C3'-C2'	-7.67	94.93	102.60
19	A7	66	A	C5-C6-N1	7.67	121.53	117.70
53	B5	131	C	O4'-C1'-N1	7.67	114.34	108.20
53	B5	1177	G	N1-C6-O6	7.67	124.50	119.90
53	B5	1777	U	O4'-C1'-N1	7.67	114.33	108.20
53	B5	2589	G	N1-C6-O6	7.67	124.50	119.90
53	B5	2677	G	O4'-C1'-N9	7.67	114.33	108.20
1	AA	1465	C	O4'-C1'-N1	7.67	114.33	108.20
1	AA	1586	G	N1-C6-O6	7.67	124.50	119.90
1	AA	1673	C	O4'-C1'-N1	7.67	114.33	108.20
53	B5	560	G	N1-C6-O6	7.67	124.50	119.90
53	B5	1187	C	O4'-C1'-N1	7.67	114.33	108.20
53	B5	2507	C	O4'-C1'-N1	7.67	114.33	108.20
53	B5	59	G	N1-C6-O6	7.67	124.50	119.90
53	B5	2794	G	N1-C6-O6	7.66	124.50	119.90
1	AA	502	U	O4'-C1'-N1	7.66	114.33	108.20
53	B5	1639	C	O4'-C1'-N1	7.66	114.33	108.20
1	AA	107	C	O4'-C1'-N1	7.66	114.33	108.20
52	B4	120	C	O4'-C1'-N1	7.66	114.33	108.20
53	B5	340	C	O4'-C1'-N1	7.66	114.33	108.20
53	B5	407	A	O4'-C1'-N9	7.66	114.33	108.20
53	B5	462	C	O4'-C1'-N1	7.66	114.33	108.20
53	B5	737	G	N1-C6-O6	7.66	124.50	119.90
53	B5	3318	G	N1-C6-O6	7.66	124.50	119.90
1	AA	519	C	O4'-C1'-N1	7.66	114.33	108.20
1	AA	1507	C	O4'-C1'-N1	7.66	114.33	108.20
53	B5	452	G	N1-C6-O6	7.66	124.49	119.90
1	AA	139	C	O4'-C1'-N1	7.66	114.33	108.20
1	AA	414	C	O4'-C1'-N1	7.66	114.33	108.20
1	AA	1594	C	O4'-C1'-N1	7.66	114.32	108.20
1	AA	1786	G	N1-C6-O6	7.66	124.49	119.90
53	B5	599	C	O4'-C1'-N1	7.66	114.32	108.20
1	AA	1566	C	O4'-C1'-N1	7.65	114.32	108.20
53	B5	2318	U	O4'-C1'-N1	7.65	114.32	108.20
53	B5	1382	G	N1-C6-O6	7.65	124.49	119.90
53	B5	2128	C	O4'-C1'-N1	7.65	114.32	108.20
53	B5	3113	A	O4'-C1'-N9	7.65	114.32	108.20
53	B5	2174	G	N1-C6-O6	7.65	124.49	119.90
53	B5	712	G	N1-C6-O6	7.65	124.49	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	2831	G	C5-C6-O6	-7.65	124.01	128.60
53	B5	3378	C	O4'-C1'-N1	7.65	114.32	108.20
1	AA	574	G	N1-C6-O6	7.64	124.49	119.90
53	B5	568	G	N1-C6-O6	7.64	124.49	119.90
53	B5	2756	C	O4'-C1'-N1	7.64	114.31	108.20
1	AA	241	U	O4'-C1'-N1	7.64	114.31	108.20
52	B4	83	C	O4'-C1'-N1	7.64	114.31	108.20
53	B5	1919	G	C5-C6-O6	-7.64	124.02	128.60
53	B5	2264	U	O4'-C1'-N1	7.64	114.31	108.20
53	B5	2732	G	N1-C6-O6	7.64	124.48	119.90
53	B5	1312	C	O4'-C1'-N1	7.64	114.31	108.20
1	AA	1200	A	O4'-C1'-N9	7.64	114.31	108.20
53	B5	166	C	O4'-C1'-N1	7.64	114.31	108.20
53	B5	297	G	N1-C6-O6	7.64	124.48	119.90
1	AA	1529	G	N1-C6-O6	7.63	124.48	119.90
53	B5	503	C	O4'-C1'-N1	7.63	114.31	108.20
53	B5	1714	A	O4'-C1'-N9	7.63	114.31	108.20
53	B5	3349	C	O4'-C1'-N1	7.63	114.31	108.20
53	B5	992	G	N1-C6-O6	7.63	124.48	119.90
1	AA	1276	C	O4'-C1'-N1	7.63	114.31	108.20
53	B5	2431	C	O4'-C1'-N1	7.63	114.31	108.20
1	AA	880	C	O4'-C1'-N1	7.63	114.30	108.20
52	B4	27	U	O4'-C1'-N1	7.63	114.30	108.20
53	B5	197	G	N1-C6-O6	7.63	124.48	119.90
53	B5	1146	C	O4'-C1'-N1	7.63	114.30	108.20
53	B5	2400	G	N1-C6-O6	7.63	124.48	119.90
1	AA	897	C	O4'-C1'-N1	7.63	114.30	108.20
53	B5	2416	U	O4'-C1'-N1	7.63	114.30	108.20
53	B5	2531	C	O4'-C1'-N1	7.63	114.30	108.20
1	AA	1644	C	O4'-C1'-N1	7.62	114.30	108.20
52	B4	36	G	N1-C6-O6	7.62	124.47	119.90
53	B5	1137	C	O4'-C1'-N1	7.62	114.30	108.20
1	AA	305	C	O4'-C1'-N1	7.62	114.30	108.20
1	AA	628	G	N1-C6-O6	7.62	124.47	119.90
53	B5	2922	G	N1-C6-O6	7.62	124.47	119.90
1	AA	846	G	N1-C6-O6	7.62	124.47	119.90
1	AA	1390	C	O4'-C1'-N1	7.62	114.30	108.20
53	B5	2714	G	N1-C6-O6	7.62	124.47	119.90
52	B4	4	C	O4'-C1'-N1	7.62	114.30	108.20
53	B5	2677	G	N1-C6-O6	7.62	124.47	119.90
53	B5	1298	C	O4'-C1'-N1	7.62	114.29	108.20
1	AA	1158	C	O4'-C1'-N1	7.62	114.29	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1546	G	C4'-C3'-C2'	7.62	110.22	102.60
53	B5	800	G	N1-C6-O6	7.62	124.47	119.90
53	B5	2199	G	N1-C6-O6	7.62	124.47	119.90
53	B5	3133	C	O4'-C1'-N1	7.61	114.29	108.20
53	B5	1857	C	O4'-C1'-N1	7.61	114.29	108.20
1	AA	1273	G	N1-C6-O6	7.61	124.47	119.90
1	AA	1603	G	O4'-C1'-N9	7.61	114.29	108.20
53	B5	3117	C	O4'-C1'-N1	7.61	114.29	108.20
1	AA	1406	G	N1-C6-O6	7.61	124.47	119.90
1	AA	683	C	O4'-C1'-N1	7.61	114.28	108.20
1	AA	1684	C	O4'-C1'-N1	7.61	114.29	108.20
23	B8	43	LYS	CB-CG-CD	7.61	131.38	111.60
53	B5	2475	G	N1-C6-O6	7.61	124.47	119.90
1	AA	1386	C	C2-N1-C1'	7.61	127.17	118.80
53	B5	2405	C	O4'-C1'-N1	7.61	114.28	108.20
53	B5	2884	C	O4'-C1'-N1	7.60	114.28	108.20
1	AA	1159	U	O4'-C1'-N1	7.60	114.28	108.20
53	B5	1898	G	O4'-C1'-N9	7.60	114.28	108.20
1	AA	1460	G	N1-C6-O6	7.60	124.46	119.90
1	AA	1589	C	O4'-C1'-N1	7.60	114.28	108.20
53	B5	2711	C	O4'-C1'-N1	7.60	114.28	108.20
53	B5	823	C	O4'-C1'-N1	7.60	114.28	108.20
1	AA	17	C	O4'-C1'-N1	7.59	114.28	108.20
1	AA	343	C	O4'-C1'-N1	7.59	114.28	108.20
1	AA	1331	U	O4'-C1'-N1	7.59	114.27	108.20
51	B3	21	G	N1-C6-O6	7.59	124.45	119.90
53	B5	2180	G	N1-C6-O6	7.59	124.45	119.90
53	B5	2918	G	P-O3'-C3'	7.59	128.81	119.70
1	AA	1170	C	O4'-C1'-N1	7.59	114.27	108.20
1	AA	1355	G	N1-C6-O6	7.59	124.45	119.90
53	B5	330	G	N1-C6-O6	7.59	124.45	119.90
53	B5	864	G	N1-C6-O6	7.59	124.45	119.90
53	B5	2751	G	N1-C6-O6	7.59	124.45	119.90
53	B5	3358	U	O4'-C1'-N1	7.59	114.27	108.20
53	B5	2273	G	N1-C6-O6	7.59	124.45	119.90
1	AA	263	C	O4'-C1'-N1	7.59	114.27	108.20
1	AA	849	C	O4'-C1'-N1	7.59	114.27	108.20
51	B3	107	G	N1-C6-O6	7.59	124.45	119.90
1	AA	1099	G	N1-C6-O6	7.58	124.45	119.90
52	B4	30	C	O4'-C1'-N1	7.58	114.27	108.20
53	B5	994	G	N1-C6-O6	7.58	124.45	119.90
53	B5	2960	C	O4'-C1'-N1	7.58	114.27	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1731	C	O4'-C1'-N1	7.58	114.27	108.20
53	B5	28	C	O4'-C1'-N1	7.58	114.27	108.20
53	B5	2605	G	N1-C6-O6	7.58	124.45	119.90
53	B5	2973	G	N1-C6-O6	7.58	124.45	119.90
53	B5	2305	G	C5-C6-O6	-7.58	124.05	128.60
1	AA	596	C	O4'-C1'-N1	7.58	114.26	108.20
53	B5	1413	G	N1-C6-O6	7.58	124.45	119.90
1	AA	543	C	O4'-C1'-N1	7.58	114.26	108.20
1	AA	1013	G	N1-C6-O6	7.58	124.45	119.90
19	A7	70	C	C5'-C4'-O4'	7.58	118.19	109.10
53	B5	2593	A	N1-C6-N6	7.58	123.14	118.60
51	B3	6	C	O4'-C1'-N1	7.57	114.26	108.20
53	B5	251	G	N1-C6-O6	7.57	124.44	119.90
53	B5	1736	G	N1-C6-O6	7.57	124.44	119.90
1	AA	949	C	O4'-C1'-N1	7.57	114.26	108.20
53	B5	1372	C	O4'-C1'-N1	7.57	114.26	108.20
53	B5	2302	G	N1-C6-O6	7.57	124.44	119.90
1	AA	686	C	O4'-C1'-N1	7.57	114.25	108.20
1	AA	1534	G	N1-C6-O6	7.57	124.44	119.90
1	AA	1580	U	O4'-C1'-N1	7.57	114.25	108.20
53	B5	360	G	N1-C6-O6	7.57	124.44	119.90
53	B5	485	A	O4'-C1'-N9	7.57	114.25	108.20
53	B5	1830	G	N1-C6-O6	7.57	124.44	119.90
53	B5	283	G	N1-C6-O6	7.57	124.44	119.90
53	B5	94	G	C5-C6-O6	-7.56	124.06	128.60
53	B5	1604	G	N1-C6-O6	7.56	124.44	119.90
1	AA	700	C	O4'-C1'-N1	7.56	114.25	108.20
1	AA	798	C	OP1-P-OP2	-7.56	108.26	119.60
53	B5	1519	G	N1-C6-O6	7.56	124.44	119.90
1	AA	1459	C	O4'-C1'-N1	7.56	114.25	108.20
1	AA	1550	U	OP1-P-OP2	-7.56	108.27	119.60
1	AA	1621	C	O4'-C1'-N1	7.56	114.25	108.20
53	B5	2719	U	O4'-C1'-N1	7.56	114.25	108.20
53	B5	2881	C	O4'-C1'-N1	7.56	114.25	108.20
53	B5	3379	C	O4'-C1'-N1	7.56	114.25	108.20
19	A7	52	U	C2-N3-C4	-7.55	122.47	127.00
53	B5	3172	G	C5-C6-O6	-7.55	124.07	128.60
1	AA	49	C	O4'-C1'-N1	7.55	114.24	108.20
1	AA	952	G	N1-C6-O6	7.55	124.43	119.90
53	B5	1156	C	O4'-C1'-N1	7.55	114.24	108.20
53	B5	1525	G	C5-C6-O6	-7.55	124.07	128.60
53	B5	1863	G	C5-C6-O6	-7.55	124.07	128.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	89	G	N1-C6-O6	7.55	124.43	119.90
1	AA	1351	G	O4'-C1'-N9	7.55	114.24	108.20
53	B5	604	G	N1-C6-O6	7.55	124.43	119.90
53	B5	180	C	O4'-C1'-N1	7.55	114.24	108.20
53	B5	1411	C	O4'-C1'-N1	7.55	114.24	108.20
1	AA	224	C	O4'-C1'-N1	7.55	114.24	108.20
53	B5	770	G	N1-C6-O6	7.55	124.43	119.90
53	B5	2898	G	N1-C6-O6	7.55	124.43	119.90
1	AA	307	G	N1-C6-O6	7.54	124.43	119.90
53	B5	1017	C	O4'-C1'-N1	7.54	114.24	108.20
53	B5	3130	A	O4'-C1'-N9	7.54	114.24	108.20
53	B5	3365	U	O4'-C1'-N1	7.54	114.24	108.20
1	AA	568	G	C5-C6-O6	-7.54	124.07	128.60
1	AA	1625	U	O4'-C1'-N1	7.54	114.23	108.20
51	B3	63	G	N1-C6-O6	7.54	124.43	119.90
53	B5	1480	G	N1-C6-O6	7.54	124.43	119.90
53	B5	2203	U	O4'-C1'-N1	7.54	114.23	108.20
53	B5	2653	C	O4'-C1'-N1	7.54	114.23	108.20
53	B5	2972	G	N1-C6-O6	7.54	124.42	119.90
1	AA	731	C	O4'-C1'-N1	7.54	114.23	108.20
53	B5	1152	G	N1-C6-O6	7.54	124.42	119.90
53	B5	1555	U	O4'-C1'-N1	7.54	114.23	108.20
53	B5	3009	G	N1-C6-O6	7.54	124.42	119.90
53	B5	893	C	O4'-C1'-N1	7.54	114.23	108.20
1	AA	365	G	N1-C6-O6	7.54	124.42	119.90
1	AA	1588	G	N1-C6-O6	7.54	124.42	119.90
53	B5	2371	G	N1-C6-O6	7.54	124.42	119.90
53	B5	2770	G	N1-C6-O6	7.54	124.42	119.90
53	B5	3329	U	O4'-C1'-N1	7.54	114.23	108.20
1	AA	325	G	O4'-C1'-N9	7.54	114.23	108.20
1	AA	595	G	N1-C6-O6	7.53	124.42	119.90
53	B5	2346	C	O4'-C1'-N1	7.53	114.23	108.20
1	AA	235	G	N1-C6-O6	7.53	124.42	119.90
1	AA	1320	C	O4'-C1'-N1	7.53	114.23	108.20
53	B5	1010	G	N1-C6-O6	7.53	124.42	119.90
53	B5	1171	G	N1-C6-O6	7.53	124.42	119.90
53	B5	1550	C	O4'-C1'-N1	7.53	114.22	108.20
1	AA	1604	C	O4'-C1'-N1	7.53	114.22	108.20
1	AA	1662	C	O4'-C1'-N1	7.53	114.22	108.20
1	AA	736	C	O4'-C1'-N1	7.53	114.22	108.20
53	B5	320	G	O4'-C1'-N9	7.53	114.22	108.20
53	B5	916	G	N1-C6-O6	7.53	124.42	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	949	C	O4'-C1'-N1	7.53	114.22	108.20
19	A7	12	U	N1-C2-N3	7.53	119.42	114.90
53	B5	31	C	O4'-C1'-N1	7.53	114.22	108.20
53	B5	212	G	N1-C6-O6	7.53	124.42	119.90
53	B5	544	C	O4'-C1'-N1	7.53	114.22	108.20
53	B5	1293	U	O4'-C1'-N1	7.53	114.22	108.20
53	B5	224	C	O4'-C1'-N1	7.52	114.22	108.20
53	B5	2284	C	O4'-C1'-N1	7.52	114.22	108.20
1	AA	1570	G	N1-C6-O6	7.52	124.41	119.90
1	AA	394	C	O4'-C1'-N1	7.52	114.22	108.20
53	B5	1012	G	N1-C6-O6	7.52	124.41	119.90
53	B5	2786	G	N1-C6-O6	7.52	124.41	119.90
53	B5	290	G	C5-C6-O6	-7.52	124.09	128.60
1	AA	409	C	O4'-C1'-N1	7.52	114.21	108.20
1	AA	1082	G	N1-C6-O6	7.52	124.41	119.90
1	AA	761	G	N1-C6-O6	7.51	124.41	119.90
53	B5	805	G	N1-C6-O6	7.51	124.41	119.90
53	B5	2365	C	O4'-C1'-N1	7.51	114.21	108.20
1	AA	523	G	N1-C6-O6	7.51	124.41	119.90
53	B5	498	A	O4'-C1'-N9	7.51	114.21	108.20
53	B5	1327	C	O4'-C1'-N1	7.51	114.21	108.20
53	B5	1611	G	N1-C6-O6	7.51	124.41	119.90
51	B3	51	G	N1-C6-O6	7.51	124.41	119.90
53	B5	493	G	N1-C6-O6	7.51	124.41	119.90
53	B5	1711	C	O4'-C1'-N1	7.51	114.21	108.20
53	B5	2325	G	N1-C6-O6	7.51	124.41	119.90
53	B5	2798	C	O4'-C1'-N1	7.51	114.21	108.20
1	AA	1525	C	O4'-C1'-N1	7.51	114.21	108.20
52	B4	56	G	N1-C6-O6	7.51	124.41	119.90
53	B5	2238	G	N1-C6-O6	7.51	124.41	119.90
53	B5	3083	G	O4'-C1'-N9	7.51	114.21	108.20
53	B5	3136	G	N1-C6-O6	7.51	124.41	119.90
51	B3	28	C	O4'-C1'-N1	7.51	114.20	108.20
53	B5	1328	C	O4'-C1'-N1	7.50	114.20	108.20
53	B5	567	G	N1-C6-O6	7.50	124.40	119.90
53	B5	3249	C	O4'-C1'-N1	7.50	114.20	108.20
1	AA	583	C	O4'-C1'-N1	7.50	114.20	108.20
53	B5	91	G	N1-C6-O6	7.50	124.40	119.90
53	B5	3180	U	O4'-C1'-N1	7.50	114.20	108.20
53	B5	2338	C	O4'-C1'-N1	7.50	114.20	108.20
53	B5	2784	G	N1-C6-O6	7.50	124.40	119.90
53	B5	2838	A	O4'-C1'-N9	7.50	114.20	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1766	G	N1-C6-O6	7.50	124.40	119.90
53	B5	335	G	N1-C6-O6	7.50	124.40	119.90
53	B5	3085	G	N1-C6-O6	7.50	124.40	119.90
1	AA	948	C	O4'-C1'-N1	7.50	114.20	108.20
51	B3	98	G	N1-C6-O6	7.50	124.40	119.90
53	B5	1513	G	N1-C6-O6	7.50	124.40	119.90
53	B5	2206	G	N1-C6-O6	7.50	124.40	119.90
53	B5	2564	G	N1-C6-O6	7.50	124.40	119.90
1	AA	860	U	C3'-C2'-C1'	-7.49	95.50	101.50
1	AA	1174	C	O4'-C1'-N1	7.49	114.19	108.20
53	B5	1815	U	O4'-C1'-N1	7.49	114.19	108.20
53	B5	2687	G	N1-C6-O6	7.49	124.39	119.90
1	AA	1202	U	O4'-C1'-N1	7.49	114.19	108.20
53	B5	1362	G	N1-C6-O6	7.49	124.39	119.90
1	AA	779	U	O4'-C1'-N1	7.49	114.19	108.20
1	AA	1270	C	O4'-C1'-N1	7.49	114.19	108.20
1	AA	1583	U	P-O3'-C3'	7.49	128.69	119.70
53	B5	138	U	O4'-C1'-N1	7.49	114.19	108.20
53	B5	741	U	O4'-C1'-N1	7.49	114.19	108.20
1	AA	1449	C	O4'-C1'-N1	7.49	114.19	108.20
53	B5	213	G	N1-C6-O6	7.49	124.39	119.90
53	B5	276	U	O4'-C1'-N1	7.49	114.19	108.20
1	AA	936	C	O4'-C1'-N1	7.48	114.19	108.20
19	A7	71	G	C4-C5-C6	-7.48	114.31	118.80
1	AA	641	G	N1-C6-O6	7.48	124.39	119.90
53	B5	33	G	N1-C6-O6	7.48	124.39	119.90
1	AA	228	G	N1-C6-O6	7.48	124.39	119.90
53	B5	1224	C	O4'-C1'-N1	7.48	114.18	108.20
53	B5	2323	G	C5-C6-O6	-7.48	124.11	128.60
53	B5	1751	G	N1-C6-O6	7.48	124.39	119.90
52	B4	135	G	N1-C6-O6	7.48	124.39	119.90
53	B5	663	C	O4'-C1'-N1	7.48	114.18	108.20
53	B5	1738	C	O4'-C1'-N1	7.48	114.18	108.20
51	B3	47	C	O4'-C1'-N1	7.48	114.18	108.20
1	AA	1539	G	N1-C6-O6	7.47	124.39	119.90
53	B5	479	U	O4'-C1'-N1	7.47	114.18	108.20
53	B5	867	G	N1-C6-O6	7.47	124.39	119.90
53	B5	3052	G	C5-C6-O6	-7.47	124.11	128.60
53	B5	1072	G	N1-C6-O6	7.47	124.38	119.90
1	AA	495	C	O4'-C1'-N1	7.47	114.18	108.20
1	AA	503	G	N1-C6-O6	7.47	124.38	119.90
53	B5	2621	G	C5-C6-O6	-7.47	124.12	128.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	2764	C	O4'-C1'-N1	7.47	114.18	108.20
1	AA	270	C	O4'-C1'-N1	7.47	114.17	108.20
53	B5	3062	G	C5-C6-O6	-7.47	124.12	128.60
1	AA	101	U	O4'-C1'-N1	7.46	114.17	108.20
1	AA	1177	C	O4'-C1'-N1	7.46	114.17	108.20
53	B5	1778	G	N1-C6-O6	7.46	124.38	119.90
1	AA	1325	G	C5-C6-O6	-7.46	124.12	128.60
53	B5	1640	G	O4'-C1'-N9	7.46	114.17	108.20
53	B5	2415	C	O4'-C1'-N1	7.46	114.17	108.20
1	AA	377	G	N1-C6-O6	7.46	124.38	119.90
53	B5	3201	C	O4'-C1'-N1	7.46	114.17	108.20
53	B5	3232	G	O4'-C1'-N9	7.46	114.17	108.20
53	B5	102	C	O4'-C1'-N1	7.46	114.17	108.20
1	AA	166	C	O4'-C1'-N1	7.46	114.17	108.20
1	AA	569	C	O4'-C1'-N1	7.46	114.17	108.20
53	B5	916	G	O4'-C1'-N9	7.46	114.17	108.20
1	AA	1564	U	O4'-C1'-N1	7.46	114.17	108.20
1	AA	455	C	O4'-C1'-N1	7.45	114.16	108.20
1	AA	1620	G	N1-C6-O6	7.45	124.37	119.90
53	B5	1706	C	O4'-C1'-N1	7.45	114.16	108.20
53	B5	1836	C	O4'-C1'-N1	7.45	114.16	108.20
53	B5	1931	U	O4'-C1'-N1	7.45	114.16	108.20
53	B5	2393	G	N1-C6-O6	7.45	124.37	119.90
53	B5	30	G	N1-C6-O6	7.45	124.37	119.90
1	AA	433	C	O4'-C1'-N1	7.45	114.16	108.20
53	B5	716	G	N1-C6-O6	7.45	124.37	119.90
53	B5	1118	C	O4'-C1'-N1	7.45	114.16	108.20
53	B5	606	C	O4'-C1'-N1	7.45	114.16	108.20
1	AA	1123	G	N1-C6-O6	7.45	124.37	119.90
53	B5	3357	U	O4'-C1'-N1	7.45	114.16	108.20
52	B4	71	G	N1-C6-O6	7.45	124.37	119.90
53	B5	1543	G	N1-C6-O6	7.45	124.37	119.90
1	AA	1135	A	N1-C6-N6	7.44	123.07	118.60
53	B5	2425	G	N1-C6-O6	7.44	124.36	119.90
53	B5	2993	G	N1-C6-O6	7.44	124.36	119.90
53	B5	3254	G	O4'-C1'-N9	7.44	114.15	108.20
1	AA	303	U	O4'-C1'-N1	7.44	114.15	108.20
53	B5	701	G	N1-C6-O6	7.44	124.36	119.90
53	B5	1635	G	N1-C6-O6	7.44	124.36	119.90
52	B4	12	A	O4'-C1'-N9	7.44	114.15	108.20
53	B5	115	A	O4'-C1'-N9	7.44	114.15	108.20
53	B5	1671	C	O4'-C1'-N1	7.44	114.15	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
52	B4	49	G	N1-C6-O6	7.44	124.36	119.90
53	B5	1690	C	O4'-C1'-N1	7.44	114.15	108.20
53	B5	538	G	N1-C6-O6	7.43	124.36	119.90
53	B5	3105	U	O4'-C1'-N1	7.43	114.15	108.20
1	AA	1231	C	O4'-C1'-N1	7.43	114.15	108.20
19	A7	7	U	C2-N3-C4	-7.43	122.54	127.00
53	B5	1256	G	C5-C6-O6	-7.43	124.14	128.60
53	B5	1595	U	O4'-C1'-N1	7.43	114.15	108.20
53	B5	2532	U	O4'-C1'-N1	7.43	114.15	108.20
1	AA	1131	C	O4'-C1'-N1	7.43	114.14	108.20
53	B5	2825	C	O4'-C1'-N1	7.43	114.14	108.20
53	B5	547	G	N1-C6-O6	7.43	124.36	119.90
53	B5	773	G	N1-C6-O6	7.43	124.36	119.90
53	B5	1375	G	N1-C6-O6	7.43	124.36	119.90
53	B5	1527	C	O4'-C1'-N1	7.43	114.14	108.20
52	B4	119	C	O4'-C1'-N1	7.43	114.14	108.20
53	B5	380	U	O4'-C1'-N1	7.43	114.14	108.20
1	AA	1612	A	C5'-C4'-O4'	7.43	118.01	109.10
52	B4	107	G	N1-C6-O6	7.43	124.36	119.90
53	B5	2552	C	O4'-C1'-N1	7.43	114.14	108.20
53	B5	3390	G	O4'-C1'-N9	7.43	114.14	108.20
1	AA	1431	U	O4'-C1'-N1	7.42	114.14	108.20
19	A7	73	A	C4'-C3'-C2'	-7.42	95.18	102.60
53	B5	363	G	N1-C6-O6	7.42	124.35	119.90
53	B5	2375	G	N1-C6-O6	7.42	124.36	119.90
19	A7	45	G	C4-C5-N7	-7.42	107.83	110.80
53	B5	337	G	N1-C6-O6	7.42	124.35	119.90
53	B5	475	G	N1-C6-O6	7.42	124.35	119.90
1	AA	141	U	O4'-C1'-N1	7.42	114.14	108.20
1	AA	396	G	N1-C6-O6	7.42	124.35	119.90
1	AA	720	G	N1-C6-O6	7.42	124.35	119.90
53	B5	420	G	N1-C6-O6	7.42	124.35	119.90
53	B5	868	C	O4'-C1'-N1	7.42	114.14	108.20
1	AA	937	G	N1-C6-O6	7.42	124.35	119.90
52	B4	34	U	O4'-C1'-N1	7.42	114.13	108.20
53	B5	1791	C	O4'-C1'-N1	7.42	114.14	108.20
53	B5	3031	G	N1-C6-O6	7.42	124.35	119.90
53	B5	1111	U	O4'-C1'-N1	7.42	114.13	108.20
53	B5	1755	C	O4'-C1'-N1	7.42	114.13	108.20
53	B5	732	C	O4'-C1'-N1	7.42	114.13	108.20
53	B5	2639	G	N1-C6-O6	7.42	124.35	119.90
1	AA	358	U	O4'-C1'-N1	7.41	114.13	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	1811	G	N1-C6-O6	7.41	124.35	119.90
53	B5	2457	G	N1-C6-O6	7.41	124.35	119.90
1	AA	799	A	O5'-P-OP1	-7.41	99.03	105.70
53	B5	2414	G	N1-C6-O6	7.41	124.35	119.90
53	B5	177	U	O4'-C1'-N1	7.41	114.13	108.20
53	B5	353	G	N1-C6-O6	7.41	124.35	119.90
53	B5	1272	C	O4'-C1'-N1	7.41	114.13	108.20
53	B5	2381	G	N1-C6-O6	7.41	124.35	119.90
53	B5	1035	G	C5-C6-O6	-7.41	124.16	128.60
53	B5	1451	C	O4'-C1'-N1	7.41	114.13	108.20
53	B5	2335	G	O4'-C1'-N9	7.41	114.13	108.20
1	AA	909	C	O4'-C1'-N1	7.41	114.12	108.20
1	AA	1709	C	O4'-C1'-N1	7.41	114.12	108.20
52	B4	139	U	O4'-C1'-N1	7.41	114.12	108.20
53	B5	1346	G	C5-C6-O6	-7.41	124.16	128.60
53	B5	2109	U	O4'-C1'-N1	7.41	114.12	108.20
53	B5	2487	U	O4'-C1'-N1	7.41	114.12	108.20
53	B5	3018	C	O4'-C1'-N1	7.41	114.12	108.20
1	AA	536	C	O4'-C1'-N1	7.40	114.12	108.20
1	AA	1379	A	O4'-C1'-N9	7.40	114.12	108.20
52	B4	32	C	O4'-C1'-N1	7.40	114.12	108.20
53	B5	150	A	O4'-C1'-N9	7.40	114.12	108.20
53	B5	424	G	N1-C6-O6	7.40	124.34	119.90
53	B5	476	G	N1-C6-O6	7.40	124.34	119.90
53	B5	1678	G	N1-C6-O6	7.40	124.34	119.90
1	AA	1781	C	O4'-C1'-N1	7.40	114.12	108.20
53	B5	566	G	N1-C6-O6	7.40	124.34	119.90
1	AA	1778	G	N1-C6-O6	7.40	124.34	119.90
53	B5	434	U	O4'-C1'-N1	7.40	114.12	108.20
53	B5	1923	C	O4'-C1'-N1	7.39	114.11	108.20
52	B4	46	G	N1-C6-O6	7.39	124.33	119.90
53	B5	270	U	O4'-C1'-N1	7.39	114.11	108.20
1	AA	691	C	O4'-C1'-N1	7.39	114.11	108.20
51	B3	65	G	N1-C6-O6	7.39	124.33	119.90
53	B5	112	U	O4'-C1'-N1	7.39	114.11	108.20
53	B5	1444	G	N1-C6-O6	7.39	124.33	119.90
53	B5	1493	G	N1-C6-O6	7.39	124.33	119.90
53	B5	110	G	O4'-C1'-N9	7.39	114.11	108.20
53	B5	2435	G	N1-C6-O6	7.39	124.33	119.90
53	B5	3118	C	O4'-C1'-N1	7.39	114.11	108.20
1	AA	1521	G	C8-N9-C1'	-7.39	117.40	127.00
19	A7	43	G	C4'-C3'-C2'	-7.39	95.21	102.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	2354	C	O4'-C1'-N1	7.39	114.11	108.20
53	B5	287	G	N1-C6-O6	7.39	124.33	119.90
53	B5	2686	A	O4'-C1'-N9	7.39	114.11	108.20
52	B4	16	G	N1-C6-O6	7.38	124.33	119.90
53	B5	679	U	O4'-C1'-N1	7.38	114.11	108.20
53	B5	2675	C	O4'-C1'-N1	7.38	114.11	108.20
53	B5	3001	C	O4'-C1'-N1	7.38	114.11	108.20
1	AA	1444	A	O4'-C1'-N9	7.38	114.11	108.20
1	AA	870	C	O4'-C1'-N1	7.38	114.11	108.20
1	AA	1516	C	O5'-C5'-C4'	-7.38	97.68	111.70
1	AA	1579	C	O4'-C1'-N1	7.38	114.11	108.20
53	B5	1230	G	O4'-C1'-N9	7.38	114.11	108.20
53	B5	1796	G	N1-C6-O6	7.38	124.33	119.90
53	B5	2717	U	O4'-C1'-N1	7.38	114.11	108.20
53	B5	3043	C	O4'-C1'-N1	7.38	114.11	108.20
52	B4	118	C	O4'-C1'-N1	7.38	114.10	108.20
53	B5	406	G	N1-C6-O6	7.38	124.33	119.90
1	AA	448	C	O4'-C1'-N1	7.38	114.10	108.20
1	AA	741	C	O4'-C1'-N1	7.38	114.10	108.20
1	AA	1306	C	O4'-C1'-N1	7.38	114.10	108.20
1	AA	1494	U	O4'-C1'-N1	7.38	114.10	108.20
19	A7	41	U	C4'-C3'-C2'	-7.38	95.22	102.60
53	B5	774	G	N1-C6-O6	7.38	124.33	119.90
53	B5	2951	G	N1-C6-O6	7.38	124.33	119.90
53	B5	1882	G	N1-C6-O6	7.38	124.33	119.90
1	AA	1118	C	O4'-C1'-N1	7.37	114.10	108.20
53	B5	2288	G	N1-C6-O6	7.37	124.33	119.90
53	B5	2335	G	N1-C6-O6	7.37	124.32	119.90
53	B5	2821	C	O4'-C1'-N1	7.37	114.10	108.20
1	AA	150	C	O4'-C1'-N1	7.37	114.10	108.20
1	AA	208	U	O4'-C1'-N1	7.37	114.10	108.20
53	B5	497	C	O4'-C1'-N1	7.37	114.10	108.20
53	B5	2467	G	C5-C6-O6	-7.37	124.18	128.60
1	AA	1218	C	O4'-C1'-N1	7.37	114.10	108.20
1	AA	1381	A	O4'-C1'-N9	7.37	114.09	108.20
19	A7	15	G	C8-N9-C4	-7.37	103.45	106.40
53	B5	1773	C	O4'-C1'-N1	7.37	114.09	108.20
1	AA	1248	C	O4'-C1'-N1	7.37	114.09	108.20
53	B5	2527	G	N1-C6-O6	7.37	124.32	119.90
1	AA	1496	G	N1-C6-O6	7.36	124.32	119.90
53	B5	728	G	C5-C6-O6	-7.36	124.18	128.60
53	B5	1175	C	O4'-C1'-N1	7.36	114.09	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	852	C	O4'-C1'-N1	7.36	114.09	108.20
53	B5	350	C	O4'-C1'-N1	7.36	114.09	108.20
1	AA	702	G	O4'-C1'-N9	7.36	114.09	108.20
53	B5	1161	G	C5-C6-O6	-7.36	124.18	128.60
53	B5	1178	G	N1-C6-O6	7.36	124.32	119.90
53	B5	2943	G	N1-C6-O6	7.36	124.31	119.90
53	B5	3089	C	O4'-C1'-N1	7.36	114.09	108.20
53	B5	502	U	O4'-C1'-N1	7.36	114.09	108.20
1	AA	152	U	O4'-C1'-N1	7.36	114.08	108.20
1	AA	646	C	O4'-C1'-N1	7.36	114.08	108.20
52	B4	63	G	N1-C6-O6	7.36	124.31	119.90
53	B5	2737	C	O4'-C1'-N1	7.36	114.08	108.20
1	AA	381	C	O4'-C1'-N1	7.35	114.08	108.20
53	B5	2239	G	N1-C6-O6	7.35	124.31	119.90
53	B5	1314	C	O4'-C1'-N1	7.35	114.08	108.20
53	B5	3229	G	N1-C6-O6	7.35	124.31	119.90
1	AA	1002	G	N1-C6-O6	7.35	124.31	119.90
53	B5	2895	G	N1-C6-O6	7.35	124.31	119.90
1	AA	362	G	N1-C6-O6	7.35	124.31	119.90
1	AA	1173	G	N1-C6-O6	7.35	124.31	119.90
19	A7	61	C	C4-C5-C6	-7.35	113.72	117.40
53	B5	842	G	N1-C6-O6	7.35	124.31	119.90
53	B5	2830	G	N1-C6-O6	7.35	124.31	119.90
1	AA	1787	G	N1-C6-O6	7.35	124.31	119.90
53	B5	744	A	O4'-C1'-N9	7.35	114.08	108.20
1	AA	990	G	C5-C6-O6	-7.34	124.19	128.60
53	B5	448	U	O4'-C1'-N1	7.34	114.08	108.20
53	B5	1383	G	C5-C6-O6	-7.34	124.19	128.60
53	B5	2481	G	N1-C6-O6	7.34	124.31	119.90
1	AA	7	G	N1-C6-O6	7.34	124.31	119.90
51	B3	96	G	N1-C6-O6	7.34	124.31	119.90
53	B5	486	U	O4'-C1'-N1	7.34	114.07	108.20
53	B5	612	U	O4'-C1'-N1	7.34	114.07	108.20
53	B5	2778	G	N1-C6-O6	7.34	124.31	119.90
53	B5	2788	C	O4'-C1'-N1	7.34	114.07	108.20
1	AA	711	U	O4'-C1'-N1	7.34	114.07	108.20
1	AA	1206	C	O4'-C1'-N1	7.34	114.07	108.20
1	AA	1492	C	O4'-C1'-N1	7.34	114.07	108.20
53	B5	505	G	N1-C6-O6	7.34	124.30	119.90
53	B5	1086	C	O4'-C1'-N1	7.34	114.07	108.20
53	B5	1814	A	O4'-C1'-N9	7.34	114.07	108.20
51	B3	44	C	O4'-C1'-N1	7.34	114.07	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	288	C	O4'-C1'-N1	7.34	114.07	108.20
1	AA	1304	U	O4'-C1'-N1	7.33	114.07	108.20
1	AA	317	C	O4'-C1'-N1	7.33	114.06	108.20
53	B5	2235	C	O4'-C1'-N1	7.33	114.06	108.20
53	B5	2857	C	O4'-C1'-N1	7.33	114.06	108.20
1	AA	1392	G	N1-C6-O6	7.33	124.30	119.90
1	AA	1015	C	O4'-C1'-N1	7.33	114.06	108.20
51	B3	106	G	N1-C6-O6	7.33	124.30	119.90
53	B5	758	C	O4'-C1'-N1	7.33	114.06	108.20
53	B5	1609	C	O4'-C1'-N1	7.33	114.06	108.20
53	B5	1821	U	O4'-C1'-N1	7.33	114.06	108.20
53	B5	1167	U	O4'-C1'-N1	7.33	114.06	108.20
53	B5	1458	U	O4'-C1'-N1	7.33	114.06	108.20
53	B5	1744	G	N1-C6-O6	7.33	124.30	119.90
53	B5	3171	U	O4'-C1'-N1	7.33	114.06	108.20
1	AA	614	C	O4'-C1'-N1	7.33	114.06	108.20
53	B5	1633	C	O4'-C1'-N1	7.33	114.06	108.20
53	B5	2301	U	O4'-C1'-N1	7.33	114.06	108.20
53	B5	2466	G	N1-C6-O6	7.33	124.30	119.90
1	AA	1303	C	O4'-C1'-N1	7.32	114.06	108.20
19	A7	35	A	N1-C2-N3	-7.32	125.64	129.30
1	AA	723	G	N1-C6-O6	7.32	124.29	119.90
53	B5	659	G	N1-C6-O6	7.32	124.29	119.90
1	AA	934	U	O4'-C1'-N1	7.32	114.06	108.20
1	AA	1330	C	O4'-C1'-N1	7.32	114.06	108.20
52	B4	19	C	O4'-C1'-N1	7.32	114.06	108.20
53	B5	1761	C	O4'-C1'-N1	7.32	114.06	108.20
53	B5	3220	G	N1-C6-O6	7.32	124.29	119.90
53	B5	3388	C	O4'-C1'-N1	7.32	114.06	108.20
1	AA	1211	C	O4'-C1'-N1	7.32	114.06	108.20
53	B5	93	C	O4'-C1'-N1	7.32	114.06	108.20
1	AA	346	G	C5-C6-O6	-7.32	124.21	128.60
1	AA	1145	C	O4'-C1'-N1	7.32	114.05	108.20
1	AA	631	G	N1-C6-O6	7.32	124.29	119.90
1	AA	1314	C	O4'-C1'-N1	7.32	114.05	108.20
53	B5	179	C	O4'-C1'-N1	7.32	114.05	108.20
53	B5	227	G	N1-C6-O6	7.32	124.29	119.90
53	B5	310	U	O4'-C1'-N1	7.32	114.05	108.20
53	B5	1941	C	O4'-C1'-N1	7.32	114.05	108.20
53	B5	2382	G	C5-C6-O6	-7.32	124.21	128.60
53	B5	2839	G	P-O3'-C3'	7.32	128.48	119.70
1	AA	858	G	N3-C4-C5	-7.31	124.94	128.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1623	C	O4'-C1'-N1	7.31	114.05	108.20
53	B5	708	G	N1-C6-O6	7.31	124.29	119.90
53	B5	2856	G	C5-C6-O6	-7.31	124.21	128.60
53	B5	918	C	O4'-C1'-N1	7.31	114.05	108.20
1	AA	1409	G	O4'-C1'-N9	7.31	114.05	108.20
51	B3	61	U	O4'-C1'-N1	7.31	114.05	108.20
1	AA	648	G	N1-C6-O6	7.31	124.28	119.90
1	AA	1622	C	O4'-C1'-N1	7.31	114.05	108.20
53	B5	412	G	O4'-C1'-N9	7.31	114.05	108.20
53	B5	2508	U	O4'-C1'-N1	7.31	114.05	108.20
46	BV	47	TYR	CB-CG-CD1	-7.31	116.62	121.00
53	B5	1284	C	O4'-C1'-N1	7.31	114.05	108.20
53	B5	1398	U	O4'-C1'-N1	7.31	114.05	108.20
53	B5	2241	U	O4'-C1'-N1	7.31	114.05	108.20
53	B5	2440	G	N1-C6-O6	7.31	124.28	119.90
1	AA	275	C	O4'-C1'-N1	7.30	114.04	108.20
53	B5	1488	G	N1-C6-O6	7.30	124.28	119.90
1	AA	338	C	O4'-C1'-N1	7.30	114.04	108.20
1	AA	1287	U	O4'-C1'-N1	7.30	114.04	108.20
53	B5	2344	U	O4'-C1'-N1	7.30	114.04	108.20
1	AA	386	G	N1-C6-O6	7.30	124.28	119.90
1	AA	1191	C	O4'-C1'-N1	7.30	114.04	108.20
19	A7	31	A	N7-C8-N9	7.30	117.45	113.80
51	B3	78	G	N1-C6-O6	7.30	124.28	119.90
51	B3	87	U	O4'-C1'-N1	7.30	114.04	108.20
53	B5	763	G	N1-C6-O6	7.30	124.28	119.90
53	B5	1325	U	O4'-C1'-N1	7.30	114.04	108.20
53	B5	1746	U	O4'-C1'-N1	7.30	114.04	108.20
53	B5	912	G	C5-C6-O6	-7.30	124.22	128.60
53	B5	1149	G	N1-C6-O6	7.30	124.28	119.90
53	B5	1895	A	O4'-C1'-N9	7.30	114.04	108.20
53	B5	2118	C	O4'-C1'-N1	7.30	114.04	108.20
53	B5	2176	U	O4'-C1'-N1	7.30	114.04	108.20
1	AA	1627	G	C5-C6-O6	-7.30	124.22	128.60
1	AA	1758	G	N1-C6-O6	7.30	124.28	119.90
53	B5	845	G	N1-C6-O6	7.30	124.28	119.90
53	B5	1731	A	O4'-C1'-N9	7.30	114.04	108.20
53	B5	2146	C	O4'-C1'-N1	7.30	114.04	108.20
53	B5	3075	G	N1-C6-O6	7.30	124.28	119.90
1	AA	443	C	O4'-C1'-N1	7.30	114.04	108.20
24	B9	14	TYR	CB-CG-CD1	-7.30	116.62	121.00
53	B5	388	G	O4'-C1'-N9	7.29	114.04	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	382	C	O4'-C1'-N1	7.29	114.03	108.20
1	AA	851	U	O4'-C1'-N1	7.29	114.03	108.20
1	AA	930	C	O4'-C1'-N1	7.29	114.03	108.20
53	B5	2378	C	O4'-C1'-N1	7.29	114.03	108.20
1	AA	18	C	O4'-C1'-N1	7.29	114.03	108.20
1	AA	160	C	O4'-C1'-N1	7.29	114.03	108.20
52	B4	28	C	O4'-C1'-N1	7.29	114.03	108.20
53	B5	698	U	O4'-C1'-N1	7.29	114.03	108.20
53	B5	803	C	O4'-C1'-N1	7.29	114.03	108.20
53	B5	2124	G	P-O3'-C3'	7.29	128.45	119.70
53	B5	3230	G	N1-C6-O6	7.29	124.27	119.90
53	B5	548	G	O4'-C1'-N9	7.29	114.03	108.20
53	B5	1623	G	N1-C6-O6	7.29	124.27	119.90
1	AA	853	G	N1-C6-O6	7.29	124.27	119.90
1	AA	967	U	O4'-C1'-N1	7.29	114.03	108.20
19	A7	62	A	C2-N3-C4	7.29	114.24	110.60
19	A7	1	G	N3-C4-C5	-7.29	124.96	128.60
53	B5	1299	U	O4'-C1'-N1	7.29	114.03	108.20
1	AA	495	C	C2-N1-C1'	7.29	126.81	118.80
1	AA	1405	G	N1-C6-O6	7.29	124.27	119.90
53	B5	911	C	O4'-C1'-N1	7.29	114.03	108.20
53	B5	2659	G	N1-C6-O6	7.29	124.27	119.90
1	AA	469	C	O4'-C1'-N1	7.28	114.03	108.20
1	AA	504	U	O4'-C1'-N1	7.28	114.03	108.20
53	B5	86	G	N1-C6-O6	7.28	124.27	119.90
53	B5	2654	C	O4'-C1'-N1	7.28	114.03	108.20
53	B5	75	G	N1-C6-O6	7.28	124.27	119.90
53	B5	1005	G	N1-C6-O6	7.28	124.27	119.90
53	B5	1094	U	O4'-C1'-N1	7.28	114.03	108.20
53	B5	1250	G	N1-C6-O6	7.28	124.27	119.90
53	B5	2408	U	O4'-C1'-N1	7.28	114.02	108.20
1	AA	1672	C	O4'-C1'-N1	7.28	114.02	108.20
19	A7	68	U	C6-N1-C2	-7.28	116.63	121.00
53	B5	979	G	N1-C6-O6	7.28	124.27	119.90
53	B5	1541	G	N1-C6-O6	7.28	124.27	119.90
1	AA	1527	C	O4'-C1'-N1	7.28	114.02	108.20
1	AA	824	G	N1-C6-O6	7.28	124.27	119.90
53	B5	226	C	O4'-C1'-N1	7.28	114.02	108.20
53	B5	315	C	O4'-C1'-N1	7.27	114.02	108.20
53	B5	944	C	O4'-C1'-N1	7.27	114.02	108.20
1	AA	552	G	N1-C6-O6	7.27	124.26	119.90
1	AA	682	C	O4'-C1'-N1	7.27	114.02	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	1371	G	N1-C6-O6	7.27	124.26	119.90
53	B5	1443	G	N1-C6-O6	7.27	124.26	119.90
1	AA	1071	G	C5-C6-O6	-7.27	124.24	128.60
51	B3	45	A	P-O3'-C3'	7.27	128.43	119.70
53	B5	565	U	O4'-C1'-N1	7.27	114.02	108.20
53	B5	1342	C	O4'-C1'-N1	7.27	114.02	108.20
53	B5	1500	G	C5-C6-O6	-7.27	124.24	128.60
1	AA	1193	C	O4'-C1'-N1	7.27	114.02	108.20
53	B5	2889	C	O4'-C1'-N1	7.27	114.02	108.20
53	B5	3300	U	O4'-C1'-N1	7.27	114.02	108.20
1	AA	419	G	N1-C6-O6	7.27	124.26	119.90
53	B5	446	U	O4'-C1'-N1	7.27	114.01	108.20
53	B5	3385	U	O4'-C1'-N1	7.27	114.01	108.20
1	AA	1162	G	N1-C6-O6	7.27	124.26	119.90
51	B3	4	U	O4'-C1'-N1	7.27	114.01	108.20
53	B5	1300	G	N1-C6-O6	7.27	124.26	119.90
53	B5	2293	C	O4'-C1'-N1	7.27	114.01	108.20
19	A7	21	A	N1-C2-N3	-7.26	125.67	129.30
53	B5	1759	C	O4'-C1'-N1	7.26	114.01	108.20
53	B5	832	G	N1-C6-O6	7.26	124.26	119.90
1	AA	1771	C	O4'-C1'-N1	7.26	114.01	108.20
53	B5	46	U	O4'-C1'-N1	7.26	114.01	108.20
53	B5	931	C	O4'-C1'-N1	7.26	114.01	108.20
53	B5	2718	U	O4'-C1'-N1	7.26	114.01	108.20
53	B5	1313	G	N1-C6-O6	7.26	124.26	119.90
53	B5	1567	U	O4'-C1'-N1	7.26	114.01	108.20
1	AA	1070	G	C5-C6-O6	-7.26	124.25	128.60
1	AA	1650	C	O4'-C1'-N1	7.26	114.00	108.20
53	B5	1765	U	O4'-C1'-N1	7.26	114.01	108.20
53	B5	2985	C	O4'-C1'-N1	7.26	114.00	108.20
53	B5	1323	G	O4'-C1'-N9	7.25	114.00	108.20
1	AA	616	G	N1-C6-O6	7.25	124.25	119.90
1	AA	639	U	O4'-C1'-N1	7.25	114.00	108.20
53	B5	264	G	N1-C6-O6	7.25	124.25	119.90
53	B5	2359	C	O4'-C1'-N1	7.25	114.00	108.20
53	B5	2528	G	N1-C6-O6	7.25	124.25	119.90
53	B5	2894	C	O4'-C1'-N1	7.25	114.00	108.20
53	B5	2944	U	O4'-C1'-N1	7.25	114.00	108.20
1	AA	281	G	N1-C6-O6	7.25	124.25	119.90
1	AA	859	A	C3'-C2'-C1'	-7.25	95.70	101.50
19	A7	69	U	N3-C4-O4	7.25	124.48	119.40
53	B5	16	A	O4'-C1'-N9	7.25	114.00	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	443	G	N1-C6-O6	7.25	124.25	119.90
53	B5	1396	C	O4'-C1'-N1	7.25	114.00	108.20
53	B5	2170	U	O4'-C1'-N1	7.25	114.00	108.20
53	B5	2240	G	N1-C6-O6	7.25	124.25	119.90
53	B5	3253	G	O4'-C1'-N9	7.25	114.00	108.20
1	AA	1116	G	N1-C6-O6	7.25	124.25	119.90
53	B5	546	C	P-O3'-C3'	7.25	128.40	119.70
53	B5	2283	G	O4'-C1'-N9	7.25	114.00	108.20
51	B3	2	G	O4'-C1'-N9	7.25	114.00	108.20
53	B5	855	U	O4'-C1'-N1	7.25	114.00	108.20
1	AA	204	G	O4'-C1'-N9	7.25	114.00	108.20
1	AA	1596	U	O4'-C1'-N1	7.25	114.00	108.20
53	B5	2185	G	N1-C6-O6	7.25	124.25	119.90
53	B5	2563	G	N1-C6-O6	7.25	124.25	119.90
53	B5	125	C	O4'-C1'-N1	7.25	114.00	108.20
53	B5	1296	C	O4'-C1'-N1	7.25	114.00	108.20
53	B5	1404	G	N1-C6-O6	7.25	124.25	119.90
53	B5	2861	U	O4'-C1'-N1	7.25	114.00	108.20
1	AA	653	C	O4'-C1'-N1	7.24	114.00	108.20
53	B5	442	G	N1-C6-O6	7.24	124.25	119.90
53	B5	454	C	O4'-C1'-N1	7.24	114.00	108.20
53	B5	1820	U	O4'-C1'-N1	7.24	114.00	108.20
53	B5	2904	U	O4'-C1'-N1	7.24	114.00	108.20
53	B5	60	A	O4'-C1'-N9	7.24	113.99	108.20
53	B5	143	G	N1-C6-O6	7.24	124.24	119.90
53	B5	650	C	O4'-C1'-N1	7.24	113.99	108.20
53	B5	1281	G	N1-C6-O6	7.24	124.25	119.90
1	AA	269	G	N1-C6-O6	7.24	124.24	119.90
51	B3	31	U	O4'-C1'-N1	7.24	113.99	108.20
53	B5	1210	U	O4'-C1'-N1	7.24	113.99	108.20
53	B5	1410	U	O4'-C1'-N1	7.24	113.99	108.20
53	B5	676	G	C5-C6-O6	-7.24	124.26	128.60
53	B5	1496	C	C6-N1-C2	-7.24	117.41	120.30
53	B5	1899	G	C5-C6-O6	-7.24	124.26	128.60
1	AA	758	U	O4'-C1'-N1	7.23	113.99	108.20
1	AA	1298	U	O4'-C1'-N1	7.23	113.99	108.20
53	B5	104	G	N1-C6-O6	7.23	124.24	119.90
52	B4	134	G	O4'-C1'-N9	7.23	113.99	108.20
53	B5	877	C	O4'-C1'-N1	7.23	113.98	108.20
53	B5	2194	G	O4'-C1'-N9	7.23	113.99	108.20
53	B5	2486	A	O4'-C1'-N9	7.23	113.99	108.20
1	AA	654	C	O4'-C1'-N1	7.23	113.98	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	2364	G	N1-C6-O6	7.23	124.24	119.90
53	B5	2409	G	N1-C6-O6	7.23	124.24	119.90
53	B5	2471	U	O4'-C1'-N1	7.23	113.98	108.20
1	AA	453	U	O4'-C1'-N1	7.23	113.98	108.20
53	B5	374	A	P-O3'-C3'	7.23	128.37	119.70
53	B5	668	G	N1-C6-O6	7.23	124.24	119.90
53	B5	1802	C	O4'-C1'-N1	7.23	113.98	108.20
53	B5	2197	C	O4'-C1'-N1	7.23	113.98	108.20
53	B5	1206	G	N1-C6-O6	7.23	124.23	119.90
53	B5	3345	G	N1-C6-O6	7.23	124.24	119.90
1	AA	90	C	O4'-C1'-N1	7.22	113.98	108.20
1	AA	186	C	O4'-C1'-N1	7.22	113.98	108.20
53	B5	634	C	O4'-C1'-N1	7.22	113.98	108.20
53	B5	1436	U	O4'-C1'-N1	7.22	113.98	108.20
53	B5	2395	G	N1-C6-O6	7.22	124.23	119.90
53	B5	2403	G	O4'-C1'-N9	7.22	113.98	108.20
1	AA	153	G	N1-C6-O6	7.22	124.23	119.90
53	B5	1388	U	O4'-C1'-N1	7.22	113.98	108.20
53	B5	1610	G	N1-C6-O6	7.22	124.23	119.90
53	B5	2730	G	N1-C6-O6	7.22	124.23	119.90
19	A7	35	A	C6-N1-C2	7.22	122.93	118.60
53	B5	494	G	N1-C6-O6	7.22	124.23	119.90
53	B5	999	G	N1-C6-O6	7.22	124.23	119.90
1	AA	942	C	O4'-C1'-N1	7.22	113.97	108.20
53	B5	890	C	O4'-C1'-N1	7.22	113.98	108.20
53	B5	2148	U	O4'-C1'-N1	7.22	113.97	108.20
1	AA	1605	G	N1-C6-O6	7.22	124.23	119.90
53	B5	286	U	P-O3'-C3'	7.22	128.36	119.70
1	AA	875	G	N1-C6-O6	7.22	124.23	119.90
19	A7	3	G	C5-C6-O6	7.22	132.93	128.60
53	B5	815	G	C5-C6-O6	-7.22	124.27	128.60
53	B5	2551	C	O4'-C1'-N1	7.22	113.97	108.20
53	B5	2834	G	N1-C6-O6	7.22	124.23	119.90
1	AA	1262	U	O4'-C1'-N1	7.21	113.97	108.20
53	B5	2110	G	C5-C6-O6	-7.21	124.27	128.60
53	B5	2254	U	O4'-C1'-N1	7.21	113.97	108.20
53	B5	491	C	O4'-C1'-N1	7.21	113.97	108.20
53	B5	2921	U	O4'-C1'-N1	7.21	113.97	108.20
53	B5	3124	G	N1-C6-O6	7.21	124.23	119.90
1	AA	598	U	O4'-C1'-N1	7.21	113.97	108.20
1	AA	668	C	O4'-C1'-N1	7.21	113.97	108.20
1	AA	781	U	O4'-C1'-N1	7.21	113.97	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
52	B4	136	G	P-O3'-C3'	7.21	128.35	119.70
53	B5	2169	G	N1-C6-O6	7.21	124.23	119.90
53	B5	2353	G	C5-C6-O6	-7.21	124.27	128.60
53	B5	2451	G	N1-C6-O6	7.21	124.23	119.90
1	AA	1186	C	O4'-C1'-N1	7.21	113.97	108.20
53	B5	391	A	O4'-C1'-N9	7.21	113.97	108.20
53	B5	3258	U	O4'-C1'-N1	7.21	113.97	108.20
1	AA	844	A	P-O3'-C3'	7.21	128.35	119.70
1	AA	856	A	C5-C6-N6	-7.21	117.93	123.70
1	AA	1039	G	N1-C6-O6	7.21	124.22	119.90
1	AA	1430	G	O4'-C1'-N9	7.21	113.97	108.20
53	B5	500	C	O4'-C1'-N1	7.21	113.97	108.20
1	AA	237	C	O4'-C1'-N1	7.21	113.97	108.20
53	B5	2283	G	N1-C6-O6	7.21	124.22	119.90
1	AA	51	A	O4'-C1'-N9	7.21	113.96	108.20
1	AA	590	C	O4'-C1'-N1	7.21	113.96	108.20
1	AA	1005	C	O4'-C1'-N1	7.20	113.96	108.20
53	B5	1672	U	O4'-C1'-N1	7.20	113.96	108.20
53	B5	1810	A	P-O3'-C3'	7.20	128.34	119.70
53	B5	1875	G	C5-C6-O6	-7.20	124.28	128.60
53	B5	3343	G	N1-C6-O6	7.20	124.22	119.90
1	AA	584	C	O4'-C1'-N1	7.20	113.96	108.20
53	B5	715	A	O4'-C1'-N9	7.20	113.96	108.20
1	AA	1127	G	C5-C6-O6	-7.20	124.28	128.60
53	B5	136	G	N1-C6-O6	7.20	124.22	119.90
53	B5	674	G	C5-C6-O6	-7.20	124.28	128.60
53	B5	745	C	O4'-C1'-N1	7.20	113.96	108.20
53	B5	1521	G	C5-C6-O6	-7.20	124.28	128.60
53	B5	1753	G	N1-C6-O6	7.20	124.22	119.90
53	B5	1776	G	C5-C6-O6	-7.20	124.28	128.60
53	B5	2412	G	N1-C6-O6	7.20	124.22	119.90
1	AA	872	G	C5-C6-O6	-7.20	124.28	128.60
53	B5	1239	C	O4'-C1'-N1	7.20	113.96	108.20
53	B5	2200	U	O4'-C1'-N1	7.20	113.96	108.20
53	B5	813	G	N1-C6-O6	7.20	124.22	119.90
53	B5	1766	G	N1-C6-O6	7.19	124.22	119.90
1	AA	763	G	N1-C6-O6	7.19	124.22	119.90
53	B5	3341	U	O4'-C1'-N1	7.19	113.95	108.20
1	AA	1538	G	N1-C6-O6	7.19	124.21	119.90
53	B5	1152	G	O4'-C1'-N9	7.19	113.95	108.20
53	B5	1234	G	N1-C6-O6	7.19	124.21	119.90
1	AA	6	G	N1-C6-O6	7.19	124.21	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	1556	C	O4'-C1'-N1	7.19	113.95	108.20
53	B5	2132	C	O4'-C1'-N1	7.19	113.95	108.20
53	B5	299	G	N1-C6-O6	7.19	124.21	119.90
1	AA	1237	G	N1-C6-O6	7.19	124.21	119.90
1	AA	1374	U	O4'-C1'-N1	7.19	113.95	108.20
53	B5	412	G	N1-C6-O6	7.19	124.21	119.90
53	B5	579	G	N1-C6-O6	7.19	124.21	119.90
53	B5	1374	G	N1-C6-O6	7.19	124.21	119.90
1	AA	1244	C	O4'-C1'-N1	7.18	113.95	108.20
1	AA	1337	U	O4'-C1'-N1	7.18	113.95	108.20
1	AA	1545	A	OP1-P-OP2	7.18	130.38	119.60
1	AA	1647	G	N1-C6-O6	7.18	124.21	119.90
1	AA	373	G	N1-C6-O6	7.18	124.21	119.90
1	AA	1552	U	OP1-P-OP2	-7.18	108.83	119.60
53	B5	56	G	N1-C6-O6	7.18	124.21	119.90
53	B5	236	G	N1-C6-O6	7.18	124.21	119.90
53	B5	639	G	C5-C6-O6	-7.18	124.29	128.60
53	B5	2753	G	N1-C6-O6	7.18	124.21	119.90
53	B5	3291	G	O4'-C1'-N9	7.18	113.94	108.20
1	AA	16	G	N1-C6-O6	7.18	124.21	119.90
53	B5	2805	G	N1-C6-O6	7.18	124.21	119.90
51	B3	25	G	N1-C6-O6	7.18	124.21	119.90
53	B5	156	G	C5-C6-O6	-7.18	124.29	128.60
53	B5	661	G	N1-C6-O6	7.18	124.21	119.90
1	AA	1120	C	O4'-C1'-N1	7.18	113.94	108.20
19	A7	5	A	C5-N7-C8	-7.18	100.31	103.90
52	B4	15	G	C5-C6-O6	-7.18	124.29	128.60
53	B5	1088	U	O4'-C1'-N1	7.18	113.94	108.20
53	B5	1207	G	N1-C6-O6	7.18	124.21	119.90
53	B5	2115	G	N1-C6-O6	7.18	124.21	119.90
1	AA	1463	C	O4'-C1'-N1	7.17	113.94	108.20
1	AA	1648	U	O4'-C1'-N1	7.17	113.94	108.20
53	B5	514	G	N1-C6-O6	7.17	124.20	119.90
53	B5	1230	G	N1-C6-O6	7.17	124.20	119.90
53	B5	3025	C	O4'-C1'-N1	7.17	113.94	108.20
1	AA	273	G	N1-C6-O6	7.17	124.20	119.90
1	AA	1757	C	O4'-C1'-N1	7.17	113.94	108.20
51	B3	58	U	O4'-C1'-N1	7.17	113.94	108.20
51	B3	102	C	O4'-C1'-N1	7.17	113.94	108.20
53	B5	1425	U	O4'-C1'-N1	7.17	113.94	108.20
53	B5	3023	U	O4'-C1'-N1	7.17	113.94	108.20
1	AA	1277	G	C5-C6-O6	-7.17	124.30	128.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	450	G	N1-C6-O6	7.17	124.20	119.90
53	B5	1267	U	O4'-C1'-N1	7.17	113.94	108.20
1	AA	175	G	N1-C6-O6	7.17	124.20	119.90
1	AA	1305	G	N1-C6-O6	7.17	124.20	119.90
53	B5	2603	G	O4'-C1'-N9	7.17	113.94	108.20
53	B5	3173	U	O4'-C1'-N1	7.17	113.94	108.20
1	AA	1033	C	O4'-C1'-N1	7.17	113.94	108.20
1	AA	1119	G	N1-C6-O6	7.17	124.20	119.90
52	B4	72	G	N1-C6-O6	7.17	124.20	119.90
53	B5	950	G	C5-C6-O6	-7.17	124.30	128.60
53	B5	1574	C	O4'-C1'-N1	7.17	113.94	108.20
53	B5	1718	G	C5-C6-O6	-7.17	124.30	128.60
53	B5	2832	C	O4'-C1'-N1	7.17	113.94	108.20
19	A7	19	G	C5-N7-C8	-7.17	100.72	104.30
51	B3	90	C	O4'-C1'-N1	7.17	113.93	108.20
52	B4	68	G	C5-C6-O6	-7.17	124.30	128.60
53	B5	1155	C	O4'-C1'-N1	7.17	113.93	108.20
53	B5	1940	G	N1-C6-O6	7.17	124.20	119.90
53	B5	2877	G	N1-C6-O6	7.17	124.20	119.90
53	B5	230	U	O4'-C1'-N1	7.17	113.93	108.20
53	B5	277	G	O4'-C1'-N9	7.17	113.93	108.20
53	B5	2722	U	O4'-C1'-N1	7.17	113.93	108.20
1	AA	1510	G	N1-C6-O6	7.16	124.20	119.90
53	B5	543	C	O4'-C1'-N1	7.16	113.93	108.20
53	B5	856	G	N1-C6-O6	7.16	124.20	119.90
53	B5	875	G	N1-C6-O6	7.16	124.20	119.90
53	B5	1432	C	O4'-C1'-N1	7.16	113.93	108.20
1	AA	825	U	O4'-C1'-N1	7.16	113.93	108.20
1	AA	1195	G	N1-C6-O6	7.16	124.20	119.90
52	B4	91	C	O4'-C1'-N1	7.16	113.93	108.20
53	B5	583	G	N1-C6-O6	7.16	124.20	119.90
53	B5	277	G	N1-C6-O6	7.16	124.19	119.90
53	B5	292	U	O4'-C1'-N1	7.16	113.93	108.20
53	B5	526	C	O4'-C1'-N1	7.16	113.93	108.20
53	B5	1279	C	O4'-C1'-N1	7.16	113.93	108.20
53	B5	1320	C	O4'-C1'-N1	7.16	113.93	108.20
53	B5	2757	U	O4'-C1'-N1	7.16	113.93	108.20
1	AA	857	U	P-O3'-C3'	-7.16	111.11	119.70
53	B5	1854	C	O4'-C1'-N1	7.16	113.92	108.20
1	AA	1240	A	O4'-C1'-N9	7.16	113.92	108.20
52	B4	124	G	N1-C6-O6	7.16	124.19	119.90
1	AA	403	G	N1-C6-O6	7.15	124.19	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1404	U	O4'-C1'-N1	7.15	113.92	108.20
53	B5	781	G	N1-C6-O6	7.15	124.19	119.90
53	B5	854	G	O4'-C1'-N9	7.15	113.92	108.20
53	B5	1140	G	C5-C6-O6	-7.15	124.31	128.60
1	AA	1624	U	O4'-C1'-N1	7.15	113.92	108.20
53	B5	1041	U	O4'-C1'-N1	7.15	113.92	108.20
53	B5	1226	G	N1-C6-O6	7.15	124.19	119.90
53	B5	2700	G	N1-C6-O6	7.15	124.19	119.90
53	B5	3301	U	O4'-C1'-N1	7.15	113.92	108.20
1	AA	432	G	N1-C6-O6	7.15	124.19	119.90
1	AA	302	U	O4'-C1'-N1	7.15	113.92	108.20
1	AA	1698	C	O4'-C1'-N1	7.15	113.92	108.20
53	B5	24	G	N1-C6-O6	7.15	124.19	119.90
53	B5	1338	C	O4'-C1'-N1	7.15	113.92	108.20
53	B5	2450	G	N1-C6-O6	7.15	124.19	119.90
53	B5	2827	U	O4'-C1'-N1	7.15	113.92	108.20
52	B4	134	G	N1-C6-O6	7.15	124.19	119.90
53	B5	2287	C	O4'-C1'-N1	7.15	113.92	108.20
53	B5	1269	U	O4'-C1'-N1	7.14	113.92	108.20
53	B5	1514	G	O4'-C1'-N9	7.14	113.92	108.20
53	B5	2957	G	N1-C6-O6	7.14	124.19	119.90
53	B5	1542	G	N1-C6-O6	7.14	124.19	119.90
1	AA	1261	G	N1-C6-O6	7.14	124.18	119.90
19	A7	12	U	C4-C5-C6	7.14	123.98	119.70
19	A7	13	C	C6-N1-C2	-7.14	117.44	120.30
53	B5	869	G	N1-C6-O6	7.14	124.18	119.90
53	B5	1023	C	O4'-C1'-N1	7.14	113.91	108.20
1	AA	724	C	O4'-C1'-N1	7.14	113.91	108.20
1	AA	1782	C	O4'-C1'-N1	7.14	113.91	108.20
53	B5	675	C	O4'-C1'-N1	7.14	113.91	108.20
53	B5	2516	U	O4'-C1'-N1	7.14	113.91	108.20
1	AA	1295	U	O4'-C1'-N1	7.14	113.91	108.20
53	B5	1142	G	N1-C6-O6	7.14	124.18	119.90
1	AA	427	C	O4'-C1'-N1	7.13	113.91	108.20
1	AA	1363	U	O4'-C1'-N1	7.13	113.91	108.20
1	AA	1542	U	O4'-C1'-N1	7.13	113.91	108.20
1	AA	1633	A	C1'-O4'-C4'	-7.13	104.19	109.90
1	AA	1737	C	O4'-C1'-N1	7.13	113.91	108.20
53	B5	1310	G	N1-C6-O6	7.13	124.18	119.90
53	B5	1501	U	O4'-C1'-N1	7.13	113.91	108.20
53	B5	1758	G	N1-C6-O6	7.13	124.18	119.90
53	B5	3090	U	O4'-C1'-N1	7.13	113.91	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	929	A	O4'-C1'-N9	7.13	113.91	108.20
53	B5	1336	U	O4'-C1'-N1	7.13	113.91	108.20
1	AA	415	C	O4'-C1'-N1	7.13	113.90	108.20
53	B5	2796	G	N1-C6-O6	7.13	124.18	119.90
53	B5	3308	C	O4'-C1'-N1	7.13	113.90	108.20
1	AA	1562	U	O4'-C1'-N1	7.13	113.90	108.20
51	B3	105	A	O4'-C1'-N9	7.13	113.90	108.20
52	B4	35	C	O4'-C1'-N1	7.13	113.90	108.20
53	B5	934	G	N1-C6-O6	7.13	124.18	119.90
53	B5	1822	C	O4'-C1'-N1	7.13	113.90	108.20
53	B5	2585	G	N1-C6-O6	7.13	124.18	119.90
53	B5	2624	G	O4'-C1'-N9	7.13	113.90	108.20
19	A7	62	A	C6-N1-C2	-7.13	114.32	118.60
53	B5	1912	U	O4'-C1'-N1	7.13	113.90	108.20
1	AA	463	U	O4'-C1'-N1	7.12	113.90	108.20
1	AA	1495	U	O4'-C1'-N1	7.12	113.90	108.20
53	B5	1789	G	N1-C6-O6	7.12	124.17	119.90
53	B5	21	G	O4'-C1'-N9	7.12	113.90	108.20
53	B5	474	G	N1-C6-O6	7.12	124.17	119.90
53	B5	1629	U	O4'-C1'-N1	7.12	113.90	108.20
53	B5	1686	U	O4'-C1'-N1	7.12	113.90	108.20
53	B5	2658	G	N1-C6-O6	7.12	124.17	119.90
1	AA	1341	A	O4'-C1'-N9	7.12	113.90	108.20
1	AA	1630	C	P-O3'-C3'	-7.12	111.16	119.70
53	B5	851	C	O4'-C1'-N1	7.12	113.90	108.20
1	AA	710	U	O4'-C1'-N1	7.12	113.90	108.20
1	AA	1153	C	O4'-C1'-N1	7.12	113.90	108.20
1	AA	1239	G	N1-C6-O6	7.12	124.17	119.90
1	AA	97	C	O4'-C1'-N1	7.12	113.90	108.20
51	B3	117	C	O4'-C1'-N1	7.12	113.89	108.20
53	B5	1579	C	O4'-C1'-N1	7.12	113.89	108.20
53	B5	2150	G	C5-C6-O6	-7.12	124.33	128.60
53	B5	2543	U	O4'-C1'-N1	7.12	113.89	108.20
1	AA	941	G	N1-C6-O6	7.12	124.17	119.90
1	AA	1743	G	N1-C6-O6	7.12	124.17	119.90
53	B5	743	C	O4'-C1'-N1	7.12	113.89	108.20
53	B5	1237	G	O4'-C1'-N9	7.12	113.89	108.20
53	B5	1868	G	C5-C6-O6	-7.12	124.33	128.60
53	B5	3059	G	N1-C6-O6	7.12	124.17	119.90
1	AA	246	G	N1-C6-O6	7.11	124.17	119.90
19	A7	28	C	N1-C2-O2	7.11	123.17	118.90
53	B5	2286	U	P-O3'-C3'	7.11	128.24	119.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1087	C	O4'-C1'-N1	7.11	113.89	108.20
53	B5	3074	G	N1-C6-O6	7.11	124.17	119.90
1	AA	1032	C	O4'-C1'-N1	7.11	113.89	108.20
52	B4	136	G	N1-C6-O6	7.11	124.17	119.90
53	B5	96	G	C5-C6-O6	-7.11	124.33	128.60
1	AA	577	G	N1-C6-O6	7.11	124.17	119.90
19	A7	65	G	O4'-C1'-N9	7.11	113.89	108.20
1	AA	1292	G	N1-C6-O6	7.11	124.16	119.90
53	B5	835	G	N1-C6-O6	7.11	124.16	119.90
1	AA	1356	C	O4'-C1'-N1	7.11	113.88	108.20
1	AA	1784	G	N1-C6-O6	7.11	124.16	119.90
19	A7	18	G	O4'-C4'-C3'	7.11	111.78	106.10
53	B5	50	U	O4'-C1'-N1	7.11	113.89	108.20
53	B5	239	G	N1-C6-O6	7.11	124.16	119.90
53	B5	1561	G	N1-C6-O6	7.11	124.16	119.90
53	B5	1646	G	N1-C6-O6	7.11	124.16	119.90
53	B5	2155	G	N1-C6-O6	7.11	124.16	119.90
53	B5	118	U	O4'-C1'-N1	7.10	113.88	108.20
1	AA	459	G	N1-C6-O6	7.10	124.16	119.90
53	B5	1268	G	N1-C6-O6	7.10	124.16	119.90
53	B5	2525	G	C5-C6-O6	-7.10	124.34	128.60
1	AA	287	G	N1-C6-O6	7.10	124.16	119.90
1	AA	1115	G	N1-C6-O6	7.10	124.16	119.90
1	AA	1660	G	N1-C6-O6	7.10	124.16	119.90
53	B5	2133	U	O4'-C1'-N1	7.10	113.88	108.20
1	AA	183	U	O4'-C1'-N1	7.10	113.88	108.20
51	B3	34	C	O4'-C1'-N1	7.10	113.88	108.20
53	B5	79	U	O4'-C1'-N1	7.10	113.88	108.20
53	B5	2463	G	O4'-C1'-N9	7.10	113.88	108.20
53	B5	3361	G	N1-C6-O6	7.10	124.16	119.90
1	AA	573	C	O4'-C1'-N1	7.10	113.88	108.20
1	AA	1516	C	C4'-C3'-C2'	7.10	109.70	102.60
1	AA	242	U	O4'-C1'-N1	7.09	113.88	108.20
1	AA	380	U	O4'-C1'-N1	7.09	113.88	108.20
53	B5	1476	G	O4'-C1'-N9	7.09	113.88	108.20
1	AA	838	G	N1-C6-O6	7.09	124.16	119.90
51	B3	41	G	N1-C6-O6	7.09	124.16	119.90
53	B5	516	A	O4'-C1'-N9	7.09	113.87	108.20
1	AA	1434	U	O4'-C1'-N1	7.09	113.87	108.20
1	AA	1714	C	O4'-C1'-N1	7.09	113.87	108.20
53	B5	2503	G	N1-C6-O6	7.09	124.15	119.90
1	AA	858	G	N1-C2-N3	-7.09	119.65	123.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	3298	C	O4'-C1'-N1	7.09	113.87	108.20
53	B5	724	U	O4'-C1'-N1	7.09	113.87	108.20
1	AA	227	U	O4'-C1'-N1	7.09	113.87	108.20
1	AA	935	G	N1-C6-O6	7.09	124.15	119.90
1	AA	1027	C	O4'-C1'-N1	7.09	113.87	108.20
1	AA	1785	C	O4'-C1'-N1	7.09	113.87	108.20
53	B5	1090	G	N1-C6-O6	7.09	124.15	119.90
53	B5	1832	C	O4'-C1'-N1	7.09	113.87	108.20
53	B5	2366	C	O4'-C1'-N1	7.09	113.87	108.20
1	AA	376	C	O4'-C1'-N1	7.08	113.87	108.20
1	AA	59	C	O4'-C1'-N1	7.08	113.87	108.20
1	AA	1102	C	O4'-C1'-N1	7.08	113.87	108.20
53	B5	682	U	O4'-C1'-N1	7.08	113.87	108.20
53	B5	1581	C	O4'-C1'-N1	7.08	113.87	108.20
53	B5	1784	G	N1-C6-O6	7.08	124.15	119.90
53	B5	1538	G	N1-C6-O6	7.08	124.15	119.90
53	B5	1824	U	O4'-C1'-N1	7.08	113.86	108.20
1	AA	865	A	C4'-C3'-O3'	7.08	127.16	113.00
53	B5	469	G	O4'-C1'-N9	7.08	113.86	108.20
53	B5	1831	U	O4'-C1'-N1	7.08	113.86	108.20
53	B5	2340	U	O4'-C1'-N1	7.08	113.86	108.20
53	B5	2749	G	N1-C6-O6	7.08	124.15	119.90
53	B5	776	U	O4'-C1'-N1	7.08	113.86	108.20
53	B5	1414	G	C5-C6-O6	-7.08	124.35	128.60
53	B5	1587	A	O4'-C1'-N9	7.08	113.86	108.20
53	B5	3121	U	P-O3'-C3'	7.08	128.19	119.70
1	AA	996	G	N1-C6-O6	7.08	124.15	119.90
1	AA	1350	U	O4'-C1'-N1	7.08	113.86	108.20
23	B8	18	TYR	CB-CG-CD2	-7.08	116.75	121.00
53	B5	1410	U	P-O3'-C3'	7.08	128.19	119.70
53	B5	1935	G	O4'-C1'-N9	7.08	113.86	108.20
53	B5	3140	G	N1-C6-O6	7.08	124.14	119.90
1	AA	1197	G	C5-C6-O6	-7.07	124.36	128.60
53	B5	654	C	O4'-C1'-N1	7.07	113.86	108.20
53	B5	1215	U	O4'-C1'-N1	7.07	113.86	108.20
1	AA	1104	G	C5-C6-O6	-7.07	124.36	128.60
1	AA	326	G	N1-C6-O6	7.07	124.14	119.90
53	B5	874	U	O4'-C1'-N1	7.07	113.86	108.20
53	B5	2248	C	O4'-C1'-N1	7.07	113.86	108.20
53	B5	2291	A	O4'-C1'-N9	7.07	113.86	108.20
52	B4	51	G	C5-C6-O6	-7.07	124.36	128.60
53	B5	127	G	C5-C6-O6	-7.07	124.36	128.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	1059	G	N1-C6-O6	7.07	124.14	119.90
1	AA	666	U	O4'-C1'-N1	7.07	113.85	108.20
1	AA	1196	G	N1-C6-O6	7.07	124.14	119.90
1	AA	1522	A	C3'-C2'-C1'	-7.07	95.85	101.50
53	B5	690	A	O4'-C1'-N9	7.07	113.85	108.20
53	B5	1402	C	O4'-C1'-N1	7.07	113.86	108.20
53	B5	3003	G	N1-C6-O6	7.07	124.14	119.90
1	AA	818	C	O4'-C1'-N1	7.07	113.85	108.20
53	B5	137	G	N1-C6-O6	7.07	124.14	119.90
1	AA	209	U	O4'-C1'-N1	7.06	113.85	108.20
53	B5	1335	C	O4'-C1'-N1	7.06	113.85	108.20
53	B5	3280	U	O4'-C1'-N1	7.06	113.85	108.20
1	AA	127	G	N1-C6-O6	7.06	124.14	119.90
53	B5	356	C	O4'-C1'-N1	7.06	113.85	108.20
53	B5	600	G	N1-C6-O6	7.06	124.14	119.90
1	AA	384	G	N1-C6-O6	7.06	124.14	119.90
51	B3	9	C	O4'-C1'-N1	7.06	113.85	108.20
53	B5	256	G	N1-C6-O6	7.06	124.14	119.90
1	AA	1143	G	N1-C6-O6	7.06	124.14	119.90
1	AA	1600	C	O4'-C1'-N1	7.06	113.85	108.20
53	B5	635	G	N1-C6-O6	7.06	124.14	119.90
53	B5	3265	C	O4'-C1'-N1	7.06	113.85	108.20
1	AA	1286	U	O4'-C1'-N1	7.06	113.85	108.20
19	A7	11	C	O3'-P-O5'	7.06	117.41	104.00
53	B5	473	G	N1-C6-O6	7.06	124.13	119.90
53	B5	1265	U	O4'-C1'-N1	7.06	113.84	108.20
1	AA	1617	C	O4'-C1'-N1	7.05	113.84	108.20
53	B5	1266	G	N1-C6-O6	7.05	124.13	119.90
53	B5	1548	C	O4'-C1'-N1	7.05	113.84	108.20
53	B5	1684	U	O4'-C1'-N1	7.05	113.84	108.20
19	A7	70	C	N3-C2-O2	-7.05	116.96	121.90
52	B4	84	C	O4'-C1'-N1	7.05	113.84	108.20
53	B5	269	G	N1-C6-O6	7.05	124.13	119.90
53	B5	1083	G	N1-C6-O6	7.05	124.13	119.90
53	B5	2784	G	O4'-C1'-N9	7.05	113.84	108.20
53	B5	2878	G	N1-C6-O6	7.05	124.13	119.90
52	B4	147	U	O4'-C1'-N1	7.05	113.84	108.20
53	B5	1073	U	O4'-C1'-N1	7.05	113.84	108.20
53	B5	3124	G	O4'-C1'-N9	7.05	113.84	108.20
53	B5	891	G	N1-C6-O6	7.05	124.13	119.90
53	B5	2171	G	N1-C6-O6	7.05	124.13	119.90
1	AA	31	C	O4'-C1'-N1	7.05	113.84	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	250	C	O4'-C1'-N1	7.05	113.84	108.20
53	B5	1108	U	O4'-C1'-N1	7.05	113.84	108.20
53	B5	2804	A	O4'-C1'-N9	7.05	113.84	108.20
53	B5	2984	C	O4'-C1'-N1	7.05	113.84	108.20
53	B5	1907	C	O4'-C1'-N1	7.05	113.84	108.20
1	AA	214	G	N1-C6-O6	7.04	124.13	119.90
1	AA	314	C	O4'-C1'-N1	7.04	113.84	108.20
1	AA	481	A	O4'-C1'-N9	7.04	113.84	108.20
1	AA	1521	G	C8-N9-C4	7.04	109.22	106.40
1	AA	1545	A	P-O3'-C3'	7.04	128.15	119.70
53	B5	367	A	O4'-C1'-N9	7.04	113.83	108.20
1	AA	634	G	N1-C6-O6	7.04	124.12	119.90
1	AA	648	G	O4'-C1'-N9	7.04	113.83	108.20
53	B5	1192	C	O4'-C1'-N1	7.04	113.83	108.20
53	B5	1344	G	N1-C6-O6	7.04	124.12	119.90
1	AA	13	C	O4'-C1'-N1	7.04	113.83	108.20
1	AA	1254	U	O4'-C1'-N1	7.04	113.83	108.20
53	B5	208	C	O4'-C1'-N1	7.04	113.83	108.20
30	BF	138	TYR	CB-CG-CD2	-7.04	116.78	121.00
52	B4	31	G	N1-C6-O6	7.04	124.12	119.90
1	AA	454	U	O4'-C1'-N1	7.04	113.83	108.20
1	AA	792	U	O5'-P-OP1	7.04	119.14	110.70
1	AA	424	C	O4'-C1'-N1	7.04	113.83	108.20
1	AA	1447	U	O4'-C1'-N1	7.04	113.83	108.20
19	A7	65	G	C8-N9-C4	-7.04	103.59	106.40
53	B5	1440	G	N1-C6-O6	7.04	124.12	119.90
53	B5	2648	G	N1-C6-O6	7.04	124.12	119.90
53	B5	3067	C	O4'-C1'-N1	7.04	113.83	108.20
1	AA	404	G	O4'-C1'-N9	7.03	113.83	108.20
52	B4	114	G	N1-C6-O6	7.03	124.12	119.90
53	B5	1229	G	O4'-C1'-N9	7.03	113.83	108.20
1	AA	1077	U	O4'-C1'-N1	7.03	113.83	108.20
53	B5	955	U	O4'-C1'-N1	7.03	113.83	108.20
53	B5	1510	G	N1-C6-O6	7.03	124.12	119.90
53	B5	3239	G	N1-C6-O6	7.03	124.12	119.90
1	AA	1635	C	O4'-C1'-N1	7.03	113.82	108.20
19	A7	9	A	C4-C5-N7	-7.03	107.19	110.70
19	A7	71	G	N1-C2-N3	-7.03	119.68	123.90
52	B4	157	U	O4'-C1'-N1	7.03	113.83	108.20
53	B5	655	C	O4'-C1'-N1	7.03	113.82	108.20
1	AA	681	U	O4'-C1'-N1	7.03	113.82	108.20
53	B5	245	U	O4'-C1'-N1	7.03	113.82	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	590	G	N1-C6-O6	7.03	124.12	119.90
1	AA	1021	C	O4'-C1'-N1	7.03	113.82	108.20
1	AA	1291	G	N1-C6-O6	7.03	124.12	119.90
53	B5	1769	G	N1-C6-O6	7.03	124.12	119.90
53	B5	2977	G	N1-C6-O6	7.03	124.12	119.90
1	AA	383	G	N1-C6-O6	7.03	124.12	119.90
1	AA	890	C	O4'-C1'-N1	7.03	113.82	108.20
19	A7	31	A	N1-C6-N6	-7.03	114.38	118.60
53	B5	978	C	O4'-C1'-N1	7.03	113.82	108.20
53	B5	1736	G	O4'-C1'-N9	7.03	113.82	108.20
53	B5	1866	C	O4'-C1'-N1	7.03	113.82	108.20
53	B5	2615	G	N1-C6-O6	7.03	124.11	119.90
53	B5	2990	G	N1-C6-O6	7.03	124.11	119.90
19	A7	64	A	N7-C8-N9	-7.02	110.29	113.80
53	B5	1439	U	O4'-C1'-N1	7.02	113.82	108.20
1	AA	800	U	O4'-C1'-N1	7.02	113.82	108.20
51	B3	37	G	N1-C6-O6	7.02	124.11	119.90
53	B5	871	U	O4'-C1'-N1	7.02	113.82	108.20
1	AA	158	U	O4'-C1'-N1	7.02	113.81	108.20
1	AA	1468	C	O4'-C1'-N1	7.02	113.81	108.20
1	AA	1203	C	O4'-C1'-N1	7.02	113.81	108.20
1	AA	1210	U	O4'-C1'-N1	7.02	113.81	108.20
53	B5	149	U	O4'-C1'-N1	7.02	113.81	108.20
53	B5	176	G	N1-C6-O6	7.02	124.11	119.90
19	A7	18	G	C3'-C2'-C1'	7.02	107.11	101.50
53	B5	909	G	N1-C6-O6	7.02	124.11	119.90
53	B5	2611	U	O4'-C1'-N1	7.02	113.81	108.20
1	AA	572	C	O4'-C1'-N1	7.01	113.81	108.20
53	B5	2328	U	O4'-C1'-N1	7.01	113.81	108.20
1	AA	968	C	O4'-C1'-N1	7.01	113.81	108.20
53	B5	3268	G	N1-C6-O6	7.01	124.11	119.90
1	AA	281	G	P-O3'-C3'	7.01	128.11	119.70
53	B5	1106	G	N1-C6-O6	7.01	124.11	119.90
53	B5	3053	G	O4'-C1'-N9	7.01	113.81	108.20
53	B5	3144	G	N1-C6-O6	7.01	124.11	119.90
53	B5	605	U	O4'-C1'-N1	7.01	113.81	108.20
53	B5	1939	G	N1-C6-O6	7.01	124.11	119.90
53	B5	2126	A	O4'-C1'-N9	7.01	113.81	108.20
53	B5	3256	G	N1-C6-O6	7.01	124.11	119.90
1	AA	709	C	O4'-C1'-N1	7.01	113.81	108.20
53	B5	90	C	O4'-C1'-N1	7.01	113.81	108.20
53	B5	358	G	N1-C6-O6	7.00	124.10	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	35	U	O4'-C1'-N1	7.00	113.80	108.20
1	AA	1755	G	N1-C6-O6	7.00	124.10	119.90
53	B5	754	G	N1-C6-O6	7.00	124.10	119.90
1	AA	677	G	N1-C6-O6	7.00	124.10	119.90
53	B5	834	U	O4'-C1'-N1	7.00	113.80	108.20
53	B5	1614	C	O4'-C1'-N1	7.00	113.80	108.20
1	AA	871	G	N1-C6-O6	7.00	124.10	119.90
1	AA	36	C	O4'-C1'-N1	7.00	113.80	108.20
1	AA	371	G	N1-C6-O6	7.00	124.10	119.90
1	AA	657	U	O4'-C1'-N1	7.00	113.80	108.20
1	AA	1272	U	O4'-C1'-N1	7.00	113.80	108.20
1	AA	1626	U	O4'-C1'-N1	7.00	113.80	108.20
1	AA	1636	G	O4'-C4'-C3'	7.00	111.70	106.10
53	B5	723	U	O4'-C1'-N1	7.00	113.80	108.20
53	B5	2292	U	O4'-C1'-N1	7.00	113.80	108.20
1	AA	422	G	N1-C6-O6	7.00	124.10	119.90
1	AA	488	G	N1-C6-O6	7.00	124.10	119.90
1	AA	503	G	O4'-C1'-N9	7.00	113.80	108.20
1	AA	1435	G	N1-C6-O6	7.00	124.10	119.90
38	BN	4	TYR	CB-CG-CD2	-7.00	116.80	121.00
44	BT	60	TYR	CB-CG-CD1	7.00	125.20	121.00
53	B5	3328	G	N1-C6-O6	7.00	124.10	119.90
1	AA	92	A	O4'-C1'-N9	7.00	113.80	108.20
1	AA	194	U	O4'-C1'-N1	7.00	113.80	108.20
1	AA	1654	U	O4'-C1'-N1	7.00	113.80	108.20
1	AA	1790	G	O4'-C1'-N9	7.00	113.80	108.20
53	B5	968	G	C5-C6-O6	-7.00	124.40	128.60
51	B3	106	G	O4'-C1'-N9	6.99	113.79	108.20
1	AA	151	G	N1-C6-O6	6.99	124.09	119.90
53	B5	540	U	O4'-C1'-N1	6.99	113.79	108.20
53	B5	819	U	O4'-C1'-N1	6.99	113.79	108.20
53	B5	2194	G	N1-C6-O6	6.99	124.09	119.90
53	B5	3309	G	O4'-C1'-N9	6.99	113.79	108.20
1	AA	123	G	N1-C6-O6	6.99	124.09	119.90
53	B5	2652	U	O4'-C1'-N1	6.99	113.79	108.20
1	AA	1565	U	O4'-C1'-N1	6.99	113.79	108.20
1	AA	1480	C	O4'-C1'-N1	6.98	113.79	108.20
53	B5	1843	C	O4'-C1'-N1	6.98	113.79	108.20
53	B5	2134	G	N1-C6-O6	6.98	124.09	119.90
53	B5	3274	G	C5-C6-O6	-6.98	124.41	128.60
19	A7	29	A	N1-C6-N6	-6.98	114.41	118.60
53	B5	196	G	C5-C6-O6	-6.98	124.41	128.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	575	G	N1-C6-O6	6.98	124.09	119.90
53	B5	2690	G	N1-C6-O6	6.98	124.09	119.90
1	AA	1512	U	O4'-C1'-N1	6.98	113.78	108.20
53	B5	126	U	O4'-C1'-N1	6.98	113.78	108.20
53	B5	831	G	N1-C6-O6	6.98	124.09	119.90
1	AA	624	G	N1-C6-O6	6.98	124.09	119.90
1	AA	1772	G	N1-C6-O6	6.98	124.09	119.90
53	B5	1213	G	N1-C6-O6	6.98	124.09	119.90
53	B5	2240	G	O4'-C1'-N9	6.98	113.78	108.20
53	B5	2850	G	O4'-C1'-N9	6.98	113.78	108.20
1	AA	1327	G	N1-C6-O6	6.97	124.08	119.90
53	B5	545	U	O4'-C1'-N1	6.97	113.78	108.20
53	B5	1615	C	O4'-C1'-N1	6.97	113.78	108.20
53	B5	1694	U	O4'-C1'-N1	6.97	113.78	108.20
51	B3	86	G	N1-C6-O6	6.97	124.08	119.90
53	B5	809	G	C5-C6-O6	-6.97	124.42	128.60
53	B5	826	G	C5-C6-O6	-6.97	124.42	128.60
1	AA	1383	G	N1-C6-O6	6.97	124.08	119.90
1	AA	1471	U	O4'-C1'-N1	6.97	113.78	108.20
53	B5	1807	G	O4'-C1'-N9	6.97	113.78	108.20
53	B5	2789	U	O4'-C1'-N1	6.97	113.78	108.20
53	B5	2914	G	N1-C6-O6	6.97	124.08	119.90
1	AA	589	C	O4'-C1'-N1	6.97	113.77	108.20
1	AA	1242	C	O4'-C1'-N1	6.97	113.77	108.20
53	B5	2728	G	N1-C6-O6	6.97	124.08	119.90
1	AA	1640	G	N1-C6-O6	6.97	124.08	119.90
53	B5	615	U	O4'-C1'-N1	6.97	113.77	108.20
53	B5	1768	U	O4'-C1'-N1	6.97	113.77	108.20
53	B5	394	G	C5-C6-O6	-6.96	124.42	128.60
53	B5	907	G	C5-C6-O6	-6.96	124.42	128.60
53	B5	1313	G	O4'-C1'-N9	6.96	113.77	108.20
53	B5	3116	G	C5-C6-O6	-6.96	124.42	128.60
53	B5	3303	G	N1-C6-O6	6.96	124.08	119.90
1	AA	276	C	O4'-C1'-N1	6.96	113.77	108.20
1	AA	1324	C	O4'-C1'-N1	6.96	113.77	108.20
53	B5	1898	G	N1-C6-O6	6.96	124.08	119.90
53	B5	2672	G	O4'-C1'-N9	6.96	113.77	108.20
1	AA	1547	C	O4'-C4'-C3'	-6.96	97.04	104.00
1	AA	1668	G	N1-C6-O6	6.96	124.08	119.90
53	B5	1044	U	O4'-C1'-N1	6.96	113.77	108.20
53	B5	1653	G	O4'-C1'-N9	6.96	113.77	108.20
53	B5	2619	G	C5-C6-O6	-6.96	124.42	128.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	796	A	C2'-C3'-O3'	6.96	124.84	113.70
1	AA	1470	C	O4'-C1'-N1	6.96	113.77	108.20
53	B5	678	G	C5-C6-O6	-6.96	124.42	128.60
53	B5	1367	G	N1-C6-O6	6.96	124.08	119.90
53	B5	3351	U	O4'-C1'-N1	6.96	113.77	108.20
1	AA	985	G	N1-C6-O6	6.96	124.08	119.90
1	AA	1680	U	O4'-C1'-N1	6.96	113.77	108.20
19	A7	51	G	C8-N9-C4	-6.96	103.62	106.40
53	B5	234	G	N1-C6-O6	6.96	124.08	119.90
53	B5	1377	G	C5-C6-O6	-6.96	124.42	128.60
53	B5	1448	U	O4'-C1'-N1	6.96	113.77	108.20
53	B5	1754	G	N1-C6-O6	6.96	124.08	119.90
53	B5	2655	U	O4'-C1'-N1	6.96	113.77	108.20
53	B5	3262	U	O4'-C1'-N1	6.96	113.77	108.20
1	AA	154	G	N1-C6-O6	6.96	124.07	119.90
1	AA	629	U	O4'-C1'-N1	6.96	113.77	108.20
53	B5	2585	G	O4'-C1'-N9	6.96	113.77	108.20
1	AA	1420	U	O4'-C1'-N1	6.96	113.76	108.20
53	B5	1063	G	N1-C6-O6	6.96	124.07	119.90
53	B5	2577	C	P-O3'-C3'	6.96	128.05	119.70
1	AA	499	U	O4'-C1'-N1	6.95	113.76	108.20
1	AA	560	U	O4'-C1'-N1	6.95	113.76	108.20
51	B3	112	G	N1-C6-O6	6.95	124.07	119.90
53	B5	432	G	N1-C6-O6	6.95	124.07	119.90
53	B5	700	C	O4'-C1'-N1	6.95	113.76	108.20
1	AA	329	G	N1-C6-O6	6.95	124.07	119.90
53	B5	3191	G	N1-C6-O6	6.95	124.07	119.90
1	AA	149	C	O4'-C1'-N1	6.95	113.76	108.20
1	AA	1290	U	O4'-C1'-N1	6.95	113.76	108.20
53	B5	740	G	N1-C6-O6	6.95	124.07	119.90
53	B5	1321	G	N1-C6-O6	6.95	124.07	119.90
53	B5	1745	C	O4'-C1'-N1	6.95	113.76	108.20
53	B5	2320	A	O4'-C1'-N9	6.95	113.76	108.20
1	AA	110	U	O4'-C1'-N1	6.94	113.75	108.20
53	B5	368	G	N1-C6-O6	6.94	124.07	119.90
53	B5	1680	G	N1-C6-O6	6.94	124.07	119.90
1	AA	704	C	O4'-C1'-N1	6.94	113.75	108.20
1	AA	1205	C	O4'-C1'-N1	6.94	113.75	108.20
1	AA	1618	C	O4'-C1'-N1	6.94	113.75	108.20
53	B5	2181	C	O4'-C1'-N1	6.94	113.75	108.20
1	AA	1451	G	N1-C6-O6	6.94	124.06	119.90
19	A7	38	A	C5-C6-N1	6.94	121.17	117.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
51	B3	59	G	N1-C6-O6	6.94	124.06	119.90
53	B5	124	U	O4'-C1'-N1	6.94	113.75	108.20
53	B5	2672	G	P-O3'-C3'	6.94	128.03	119.70
1	AA	798	C	C2'-C3'-O3'	6.94	124.80	113.70
1	AA	1666	G	N1-C6-O6	6.94	124.06	119.90
51	B3	76	U	O4'-C1'-N1	6.94	113.75	108.20
53	B5	2782	U	O4'-C1'-N1	6.94	113.75	108.20
19	A7	29	A	C4-C5-C6	-6.93	113.53	117.00
53	B5	785	G	C5-C6-O6	-6.93	124.44	128.60
53	B5	1013	G	N1-C6-O6	6.93	124.06	119.90
53	B5	1307	G	N1-C6-O6	6.93	124.06	119.90
53	B5	2530	G	N1-C6-O6	6.93	124.06	119.90
1	AA	229	U	O4'-C1'-N1	6.93	113.75	108.20
1	AA	1006	C	O4'-C1'-N1	6.93	113.75	108.20
53	B5	1883	A	O4'-C1'-N9	6.93	113.75	108.20
1	AA	282	C	O4'-C1'-N1	6.93	113.75	108.20
1	AA	750	U	O4'-C1'-N1	6.93	113.75	108.20
1	AA	64	U	O4'-C1'-N1	6.93	113.74	108.20
1	AA	759	U	O4'-C1'-N1	6.93	113.74	108.20
1	AA	1184	G	N1-C6-O6	6.93	124.06	119.90
1	AA	1474	C	O4'-C1'-N1	6.93	113.74	108.20
19	A7	20	G	N9-C1'-C2'	6.93	123.01	114.00
51	B3	88	G	C5-C6-O6	-6.93	124.44	128.60
53	B5	703	G	N1-C6-O6	6.93	124.06	119.90
53	B5	2920	U	O4'-C1'-N1	6.93	113.74	108.20
53	B5	850	U	O4'-C1'-N1	6.93	113.74	108.20
53	B5	1322	U	O4'-C1'-N1	6.93	113.74	108.20
53	B5	1936	A	O4'-C1'-N9	6.93	113.74	108.20
53	B5	5	G	N1-C6-O6	6.93	124.06	119.90
53	B5	2632	G	N1-C6-O6	6.93	124.06	119.90
53	B5	3153	U	O4'-C1'-N1	6.93	113.74	108.20
19	A7	41	U	N1-C2-N3	6.92	119.06	114.90
53	B5	1607	U	O4'-C1'-N1	6.92	113.74	108.20
1	AA	626	U	O4'-C1'-N1	6.92	113.74	108.20
1	AA	776	G	N1-C6-O6	6.92	124.05	119.90
52	B4	152	G	N1-C6-O6	6.92	124.05	119.90
53	B5	903	U	O4'-C1'-N1	6.92	113.74	108.20
53	B5	1363	A	O4'-C1'-N9	6.92	113.74	108.20
53	B5	1788	C	O4'-C1'-N1	6.92	113.74	108.20
53	B5	2940	A	O4'-C1'-N9	6.92	113.74	108.20
1	AA	953	G	N1-C6-O6	6.92	124.05	119.90
52	B4	140	G	N1-C6-O6	6.92	124.05	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	217	U	O4'-C1'-N1	6.92	113.74	108.20
53	B5	447	U	O4'-C1'-N1	6.92	113.74	108.20
53	B5	1138	U	O4'-C1'-N1	6.92	113.74	108.20
53	B5	3380	U	O4'-C1'-N1	6.92	113.74	108.20
1	AA	1425	G	C5-C6-O6	-6.92	124.45	128.60
53	B5	636	C	O4'-C1'-N1	6.92	113.73	108.20
53	B5	954	U	O4'-C1'-N1	6.92	113.73	108.20
53	B5	1897	G	N1-C6-O6	6.92	124.05	119.90
1	AA	146	U	O4'-C1'-N1	6.92	113.73	108.20
1	AA	297	U	O4'-C1'-N1	6.92	113.73	108.20
1	AA	1639	C	O4'-C1'-N1	6.92	113.73	108.20
53	B5	702	C	O4'-C1'-N1	6.92	113.73	108.20
53	B5	2754	G	N1-C6-O6	6.92	124.05	119.90
1	AA	94	U	O4'-C1'-N1	6.92	113.73	108.20
1	AA	243	G	N1-C6-O6	6.92	124.05	119.90
1	AA	823	G	N1-C6-O6	6.92	124.05	119.90
1	AA	1505	G	N1-C6-O6	6.92	124.05	119.90
53	B5	1618	G	N1-C6-O6	6.92	124.05	119.90
1	AA	715	U	O4'-C1'-N1	6.91	113.73	108.20
53	B5	1700	G	N1-C6-O6	6.91	124.05	119.90
53	B5	1889	G	N1-C6-O6	6.91	124.05	119.90
53	B5	2162	U	O4'-C1'-N1	6.91	113.73	108.20
53	B5	3238	G	N1-C6-O6	6.91	124.05	119.90
1	AA	1046	U	O4'-C1'-N1	6.91	113.73	108.20
1	AA	449	C	O4'-C1'-N1	6.91	113.73	108.20
1	AA	1531	C	O4'-C1'-N1	6.91	113.73	108.20
19	A7	64	A	C5-C6-N1	6.91	121.16	117.70
53	B5	52	A	C5-C6-N6	-6.91	118.17	123.70
53	B5	223	U	O4'-C1'-N1	6.91	113.73	108.20
53	B5	746	A	O4'-C1'-N9	6.91	113.73	108.20
53	B5	1078	U	O4'-C1'-N1	6.91	113.73	108.20
53	B5	833	G	N1-C6-O6	6.91	124.05	119.90
53	B5	1081	U	O4'-C1'-N1	6.91	113.73	108.20
53	B5	1092	C	O4'-C1'-N1	6.91	113.73	108.20
53	B5	1772	U	O4'-C1'-N1	6.91	113.73	108.20
53	B5	2981	U	O4'-C1'-N1	6.91	113.73	108.20
53	B5	2406	C	O4'-C1'-N1	6.91	113.72	108.20
1	AA	1243	U	O4'-C1'-N1	6.91	113.72	108.20
1	AA	1569	C	O4'-C1'-N1	6.91	113.72	108.20
53	B5	10	C	O4'-C1'-N1	6.91	113.72	108.20
53	B5	2704	A	O4'-C1'-N9	6.91	113.72	108.20
1	AA	102	U	O4'-C1'-N1	6.90	113.72	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	703	G	O4'-C1'-N9	6.90	113.72	108.20
1	AA	1007	G	N1-C6-O6	6.90	124.04	119.90
53	B5	231	G	O4'-C1'-N9	6.90	113.72	108.20
53	B5	1228	C	O4'-C1'-N1	6.90	113.72	108.20
53	B5	1775	G	N1-C6-O6	6.90	124.04	119.90
53	B5	2554	U	O4'-C1'-N1	6.90	113.72	108.20
53	B5	2937	G	N1-C6-O6	6.90	124.04	119.90
53	B5	3192	U	O4'-C1'-N1	6.90	113.72	108.20
53	B5	2396	G	N1-C6-O6	6.90	124.04	119.90
53	B5	2972	G	O4'-C1'-N9	6.90	113.72	108.20
1	AA	413	U	O4'-C1'-N1	6.90	113.72	108.20
53	B5	582	G	N1-C6-O6	6.90	124.04	119.90
53	B5	1121	U	O4'-C1'-N1	6.90	113.72	108.20
53	B5	2826	U	O4'-C1'-N1	6.90	113.72	108.20
1	AA	698	U	O4'-C1'-N1	6.89	113.72	108.20
53	B5	507	U	O4'-C1'-N1	6.89	113.72	108.20
53	B5	762	U	O4'-C1'-N1	6.89	113.72	108.20
1	AA	901	G	N1-C6-O6	6.89	124.03	119.90
53	B5	464	U	O4'-C1'-N1	6.89	113.71	108.20
53	B5	1478	C	O4'-C1'-N1	6.89	113.71	108.20
53	B5	2216	G	C5-C6-O6	-6.89	124.47	128.60
53	B5	362	U	O4'-C1'-N1	6.89	113.71	108.20
52	B4	74	U	O4'-C1'-N1	6.89	113.71	108.20
53	B5	2134	G	O4'-C1'-N9	6.89	113.71	108.20
53	B5	2998	U	O4'-C1'-N1	6.89	113.71	108.20
53	B5	3363	U	O4'-C1'-N1	6.89	113.71	108.20
53	B5	1817	G	N1-C6-O6	6.89	124.03	119.90
1	AA	1164	G	O4'-C1'-N9	6.89	113.71	108.20
1	AA	1602	U	O4'-C1'-N1	6.89	113.71	108.20
53	B5	1014	U	O4'-C1'-N1	6.89	113.71	108.20
1	AA	786	C	O4'-C1'-N1	6.88	113.71	108.20
53	B5	1127	G	C5-C6-O6	-6.88	124.47	128.60
52	B4	126	A	O4'-C1'-N9	6.88	113.71	108.20
53	B5	143	G	O4'-C1'-N9	6.88	113.71	108.20
53	B5	419	G	N1-C6-O6	6.88	124.03	119.90
53	B5	1591	G	N1-C6-O6	6.88	124.03	119.90
53	B5	756	U	O4'-C1'-N1	6.88	113.70	108.20
53	B5	892	U	O4'-C1'-N1	6.88	113.70	108.20
53	B5	1233	G	O4'-C1'-N9	6.88	113.70	108.20
53	B5	2555	G	N1-C6-O6	6.88	124.03	119.90
53	B5	3253	G	N1-C6-O6	6.88	124.03	119.90
1	AA	1472	G	C5-C6-O6	-6.88	124.47	128.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1555	U	O4'-C1'-N1	6.88	113.70	108.20
1	AA	1683	G	N1-C6-O6	6.88	124.03	119.90
53	B5	97	U	O4'-C1'-N1	6.88	113.70	108.20
53	B5	2604	U	O4'-C1'-N1	6.88	113.70	108.20
53	B5	2720	G	N1-C6-O6	6.88	124.03	119.90
53	B5	2762	A	O4'-C1'-N9	6.88	113.70	108.20
53	B5	3321	C	O4'-C1'-N1	6.88	113.70	108.20
1	AA	885	G	N1-C6-O6	6.88	124.03	119.90
1	AA	1483	C	O4'-C1'-N1	6.88	113.70	108.20
52	B4	116	G	N1-C6-O6	6.88	124.03	119.90
53	B5	512	U	P-O3'-C3'	6.88	127.95	119.70
53	B5	821	U	O4'-C1'-N1	6.88	113.70	108.20
53	B5	2791	G	N1-C6-O6	6.88	124.03	119.90
53	B5	3109	G	O4'-C1'-N9	6.88	113.70	108.20
53	B5	58	G	N1-C6-O6	6.88	124.03	119.90
53	B5	1700	G	O4'-C1'-N9	6.88	113.70	108.20
1	AA	1227	U	O4'-C1'-N1	6.87	113.70	108.20
1	AA	1264	G	N1-C6-O6	6.87	124.02	119.90
23	B8	43	LYS	CG-CD-CE	6.87	132.52	111.90
53	B5	1211	U	O4'-C1'-N1	6.87	113.70	108.20
53	B5	1427	U	O4'-C1'-N1	6.87	113.70	108.20
53	B5	1593	A	O4'-C1'-N9	6.87	113.70	108.20
1	AA	285	G	N1-C6-O6	6.87	124.02	119.90
53	B5	339	C	O4'-C1'-N1	6.87	113.70	108.20
1	AA	294	C	O4'-C1'-N1	6.87	113.70	108.20
1	AA	988	U	O4'-C1'-N1	6.87	113.69	108.20
1	AA	1478	G	N1-C6-O6	6.87	124.02	119.90
53	B5	779	G	N1-C6-O6	6.87	124.02	119.90
53	B5	786	A	O4'-C1'-N9	6.87	113.70	108.20
53	B5	1218	U	O4'-C1'-N1	6.87	113.69	108.20
53	B5	1902	G	C5-C6-O6	-6.87	124.48	128.60
53	B5	2706	G	N1-C6-O6	6.87	124.02	119.90
1	AA	612	U	O4'-C1'-N1	6.87	113.69	108.20
53	B5	546	C	O4'-C1'-N1	6.87	113.69	108.20
53	B5	798	G	N1-C6-O6	6.87	124.02	119.90
53	B5	1536	G	N1-C6-O6	6.87	124.02	119.90
53	B5	811	U	O4'-C1'-N1	6.87	113.69	108.20
53	B5	1157	G	C5-C6-O6	-6.87	124.48	128.60
53	B5	1661	G	N1-C6-O6	6.87	124.02	119.90
53	B5	2201	G	N1-C6-O6	6.87	124.02	119.90
1	AA	1438	C	O4'-C1'-N1	6.86	113.69	108.20
53	B5	261	U	O4'-C1'-N1	6.86	113.69	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
52	B4	81	U	O4'-C1'-N1	6.86	113.69	108.20
53	B5	3259	U	O4'-C1'-N1	6.86	113.69	108.20
1	AA	121	U	O4'-C1'-N1	6.86	113.69	108.20
1	AA	1384	G	N1-C6-O6	6.86	124.02	119.90
1	AA	1557	A	O4'-C1'-N9	6.86	113.69	108.20
19	A7	60	C	O4'-C1'-N1	6.86	113.69	108.20
51	B3	10	C	O4'-C1'-N1	6.86	113.69	108.20
53	B5	1795	U	P-O3'-C3'	6.86	127.93	119.70
53	B5	937	G	N1-C6-O6	6.86	124.02	119.90
53	B5	1033	U	O4'-C1'-N1	6.86	113.69	108.20
1	AA	1063	C	O4'-C1'-N1	6.86	113.69	108.20
1	AA	1228	U	O4'-C1'-N1	6.86	113.69	108.20
51	B3	11	A	O4'-C1'-N9	6.86	113.69	108.20
53	B5	27	C	O4'-C1'-N1	6.86	113.69	108.20
53	B5	1634	G	N1-C6-O6	6.86	124.01	119.90
53	B5	1790	G	C5-C6-O6	-6.86	124.49	128.60
53	B5	2647	A	O4'-C1'-N9	6.86	113.69	108.20
53	B5	3065	G	N1-C6-O6	6.86	124.02	119.90
53	B5	160	G	N1-C6-O6	6.86	124.01	119.90
53	B5	220	G	N1-C6-O6	6.86	124.01	119.90
53	B5	1000	C	O4'-C1'-N1	6.86	113.68	108.20
1	AA	787	G	O4'-C1'-N9	6.85	113.68	108.20
19	A7	70	C	O4'-C1'-N1	6.85	113.68	108.20
53	B5	2177	G	C5-C6-O6	-6.85	124.49	128.60
1	AA	1347	U	O4'-C1'-N1	6.85	113.68	108.20
51	B3	56	G	O4'-C1'-N9	6.85	113.68	108.20
53	B5	587	U	O4'-C1'-N1	6.85	113.68	108.20
53	B5	2210	G	C5-C6-O6	-6.85	124.49	128.60
13	AN	16	LYS	N-CA-C	6.85	129.49	111.00
53	B5	120	G	C5-C6-O6	-6.85	124.49	128.60
53	B5	793	C	O4'-C1'-N1	6.85	113.68	108.20
53	B5	2436	U	O4'-C1'-N1	6.85	113.68	108.20
53	B5	3255	U	O4'-C1'-N1	6.85	113.68	108.20
1	AA	279	G	N1-C6-O6	6.85	124.01	119.90
1	AA	1614	G	N1-C6-O6	6.85	124.01	119.90
53	B5	84	U	O4'-C1'-N1	6.85	113.68	108.20
53	B5	204	A	O4'-C1'-N9	6.85	113.68	108.20
53	B5	601	U	O4'-C1'-N1	6.85	113.68	108.20
53	B5	995	G	N1-C6-O6	6.85	124.01	119.90
53	B5	1650	G	N1-C6-O6	6.85	124.01	119.90
53	B5	2843	U	O4'-C1'-N1	6.85	113.68	108.20
1	AA	1462	G	N1-C6-O6	6.85	124.01	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	3026	G	N1-C6-O6	6.85	124.01	119.90
1	AA	785	U	O4'-C1'-N1	6.84	113.68	108.20
1	AA	1452	G	N1-C6-O6	6.84	124.01	119.90
19	A7	45	G	N9-C4-C5	6.84	108.14	105.40
51	B3	84	G	N1-C6-O6	6.84	124.01	119.90
53	B5	506	U	O4'-C1'-N1	6.84	113.67	108.20
53	B5	2111	G	C5-C6-O6	-6.84	124.49	128.60
53	B5	3102	G	C5-C6-O6	-6.84	124.49	128.60
1	AA	1365	G	N1-C6-O6	6.84	124.01	119.90
53	B5	1089	G	N1-C6-O6	6.84	124.00	119.90
53	B5	1126	G	N1-C6-O6	6.84	124.00	119.90
53	B5	1457	U	O4'-C1'-N1	6.84	113.67	108.20
53	B5	1710	C	O4'-C1'-N1	6.84	113.67	108.20
53	B5	1837	U	O4'-C1'-N1	6.84	113.67	108.20
53	B5	3071	U	O4'-C1'-N1	6.84	113.67	108.20
1	AA	336	G	N1-C6-O6	6.84	124.00	119.90
53	B5	1532	C	O4'-C1'-N1	6.84	113.67	108.20
53	B5	2270	A	O4'-C1'-N9	6.84	113.67	108.20
53	B5	1415	U	O4'-C1'-N1	6.84	113.67	108.20
53	B5	1894	U	O4'-C1'-N1	6.84	113.67	108.20
53	B5	2645	G	N1-C6-O6	6.84	124.00	119.90
1	AA	212	U	O4'-C1'-N1	6.84	113.67	108.20
53	B5	181	U	O4'-C1'-N1	6.84	113.67	108.20
1	AA	915	U	O4'-C1'-N1	6.83	113.67	108.20
1	AA	999	C	O4'-C1'-N1	6.83	113.67	108.20
53	B5	145	G	N1-C6-O6	6.83	124.00	119.90
53	B5	2112	U	O4'-C1'-N1	6.83	113.67	108.20
53	B5	2651	G	C5-C6-O6	-6.83	124.50	128.60
53	B5	445	G	N1-C6-O6	6.83	124.00	119.90
1	AA	1656	G	O4'-C1'-N9	6.83	113.67	108.20
53	B5	969	C	O4'-C1'-N1	6.83	113.67	108.20
53	B5	1274	A	O4'-C1'-N9	6.83	113.67	108.20
53	B5	2388	U	O4'-C1'-N1	6.83	113.67	108.20
1	AA	14	C	O4'-C1'-N1	6.83	113.66	108.20
1	AA	867	G	O4'-C4'-C3'	-6.83	97.17	104.00
1	AA	1008	U	O4'-C1'-N1	6.83	113.66	108.20
1	AA	1746	G	N1-C6-O6	6.83	124.00	119.90
53	B5	2207	A	O4'-C1'-N9	6.83	113.66	108.20
53	B5	2474	G	N1-C6-O6	6.83	124.00	119.90
1	AA	647	G	N1-C6-O6	6.83	124.00	119.90
1	AA	946	U	O4'-C1'-N1	6.83	113.66	108.20
53	B5	2506	U	O4'-C1'-N1	6.83	113.66	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	18	G	N1-C6-O6	6.83	124.00	119.90
53	B5	178	U	O4'-C1'-N1	6.83	113.66	108.20
53	B5	2322	C	O4'-C1'-N1	6.83	113.66	108.20
1	AA	842	C	O4'-C1'-N1	6.82	113.66	108.20
53	B5	1340	G	C5-C6-O6	-6.82	124.51	128.60
53	B5	2823	G	C5-C6-O6	-6.82	124.51	128.60
1	AA	1716	G	N1-C6-O6	6.82	123.99	119.90
53	B5	2454	G	N1-C6-O6	6.82	123.99	119.90
1	AA	888	U	O4'-C1'-N1	6.82	113.66	108.20
52	B4	133	G	N1-C6-O6	6.82	123.99	119.90
53	B5	1097	G	N1-C6-O6	6.82	123.99	119.90
53	B5	1748	G	C5-C6-O6	-6.82	124.51	128.60
1	AA	1051	U	O4'-C1'-N1	6.82	113.66	108.20
53	B5	1855	U	O4'-C1'-N1	6.82	113.66	108.20
53	B5	2369	G	N1-C6-O6	6.82	123.99	119.90
1	AA	120	U	O4'-C1'-N1	6.82	113.65	108.20
51	B3	1	G	C5-C6-O6	-6.82	124.51	128.60
53	B5	3231	U	O4'-C1'-N1	6.82	113.65	108.20
1	AA	185	U	O4'-C1'-N1	6.82	113.65	108.20
53	B5	1117	G	C5-C6-O6	-6.82	124.51	128.60
53	B5	2548	C	C6-N1-C2	-6.82	117.57	120.30
1	AA	1453	G	C5-C6-O6	-6.81	124.51	128.60
53	B5	838	G	N1-C6-O6	6.81	123.99	119.90
53	B5	2598	G	N1-C6-O6	6.81	123.99	119.90
53	B5	2829	U	O4'-C1'-N1	6.81	113.65	108.20
1	AA	858	G	C6-C5-N7	-6.81	126.31	130.40
53	B5	64	G	C5-C6-O6	-6.81	124.51	128.60
53	B5	2333	C	O4'-C1'-N1	6.81	113.65	108.20
53	B5	2394	G	O4'-C1'-N9	6.81	113.65	108.20
1	AA	1576	U	O4'-C1'-N1	6.81	113.65	108.20
1	AA	304	U	O4'-C1'-N1	6.81	113.65	108.20
52	B4	122	U	O4'-C1'-N1	6.81	113.65	108.20
52	B4	158	U	O4'-C1'-N1	6.81	113.65	108.20
53	B5	1370	G	N1-C6-O6	6.81	123.99	119.90
1	AA	1312	U	O4'-C1'-N1	6.81	113.65	108.20
53	B5	1833	G	C5-C6-O6	-6.81	124.52	128.60
53	B5	1844	C	O4'-C1'-N1	6.81	113.65	108.20
53	B5	2723	U	O4'-C1'-N1	6.81	113.65	108.20
53	B5	3232	G	N1-C6-O6	6.81	123.98	119.90
1	AA	468	A	O4'-C1'-N9	6.81	113.64	108.20
1	AA	1042	C	O4'-C1'-N1	6.81	113.64	108.20
1	AA	1076	U	O4'-C1'-N1	6.81	113.64	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1466	U	O4'-C1'-N1	6.81	113.64	108.20
53	B5	331	G	N1-C6-O6	6.81	123.98	119.90
53	B5	1673	G	N1-C6-O6	6.81	123.98	119.90
53	B5	2446	U	O4'-C1'-N1	6.81	113.64	108.20
53	B5	2776	C	O4'-C1'-N1	6.81	113.64	108.20
53	B5	3082	C	O4'-C1'-N1	6.81	113.64	108.20
1	AA	96	G	N1-C6-O6	6.80	123.98	119.90
53	B5	1848	G	N1-C6-O6	6.80	123.98	119.90
53	B5	2541	U	O4'-C1'-N1	6.80	113.64	108.20
53	B5	2745	G	N1-C6-O6	6.80	123.98	119.90
53	B5	2763	U	O4'-C1'-N1	6.80	113.64	108.20
1	AA	659	C	O4'-C1'-N1	6.80	113.64	108.20
51	B3	36	C	O4'-C1'-N1	6.80	113.64	108.20
51	B3	49	G	N1-C6-O6	6.80	123.98	119.90
53	B5	938	C	O4'-C1'-N1	6.80	113.64	108.20
53	B5	1242	G	N1-C6-O6	6.80	123.98	119.90
53	B5	1782	U	O4'-C1'-N1	6.80	113.64	108.20
53	B5	2581	U	O4'-C1'-N1	6.80	113.64	108.20
53	B5	3080	G	N1-C6-O6	6.80	123.98	119.90
53	B5	1186	G	C5-C6-O6	-6.80	124.52	128.60
53	B5	1417	G	N1-C6-O6	6.80	123.98	119.90
53	B5	1794	G	C5-C6-O6	-6.80	124.52	128.60
53	B5	3263	G	N1-C6-O6	6.80	123.98	119.90
53	B5	2411	U	O4'-C1'-N1	6.80	113.64	108.20
53	B5	3064	U	O4'-C1'-N1	6.80	113.64	108.20
19	A7	18	G	C2-N3-C4	-6.80	108.50	111.90
53	B5	641	C	O4'-C1'-N1	6.80	113.64	108.20
53	B5	1659	U	O4'-C1'-N1	6.80	113.64	108.20
53	B5	299	G	O4'-C1'-N9	6.79	113.64	108.20
53	B5	887	G	N1-C6-O6	6.79	123.98	119.90
19	A7	30	G	C2-N3-C4	6.79	115.30	111.90
19	A7	44	A	N1-C2-N3	-6.79	125.91	129.30
53	B5	2596	U	O4'-C1'-N1	6.79	113.63	108.20
53	B5	3100	U	O4'-C1'-N1	6.79	113.63	108.20
52	B4	55	U	O4'-C1'-N1	6.79	113.63	108.20
53	B5	1431	G	N1-C6-O6	6.79	123.97	119.90
53	B5	2983	C	O4'-C1'-N1	6.79	113.63	108.20
1	AA	447	U	O4'-C1'-N1	6.79	113.63	108.20
1	AA	575	C	O4'-C1'-N1	6.79	113.63	108.20
52	B4	100	U	O4'-C1'-N1	6.79	113.63	108.20
52	B4	123	G	N1-C6-O6	6.79	123.97	119.90
53	B5	17	G	N1-C6-O6	6.79	123.97	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	1101	G	N1-C6-O6	6.79	123.97	119.90
53	B5	2219	A	P-O3'-C3'	6.79	127.85	119.70
1	AA	361	C	O4'-C1'-N1	6.79	113.63	108.20
1	AA	1545	A	C2'-C3'-O3'	6.79	124.56	113.70
53	B5	1769	G	O4'-C1'-N9	6.79	113.63	108.20
53	B5	1808	G	N1-C6-O6	6.79	123.97	119.90
53	B5	2668	U	O4'-C1'-N1	6.79	113.63	108.20
1	AA	1454	C	O4'-C1'-N1	6.78	113.63	108.20
19	A7	75	C	C5-C4-N4	-6.78	115.45	120.20
51	B3	67	C	O4'-C1'-N1	6.78	113.63	108.20
53	B5	41	G	N1-C6-O6	6.78	123.97	119.90
53	B5	168	U	O4'-C1'-N1	6.78	113.63	108.20
53	B5	322	U	O4'-C1'-N1	6.78	113.63	108.20
52	B4	20	U	O4'-C1'-N1	6.78	113.63	108.20
53	B5	588	G	N1-C6-O6	6.78	123.97	119.90
1	AA	240	U	C2-N1-C1'	6.78	125.84	117.70
53	B5	764	U	O4'-C1'-N1	6.78	113.62	108.20
53	B5	1113	G	N1-C6-O6	6.78	123.97	119.90
53	B5	2676	A	O4'-C1'-N9	6.78	113.62	108.20
1	AA	57	G	C5-C6-O6	-6.78	124.53	128.60
1	AA	1019	A	O4'-C1'-N9	6.78	113.62	108.20
1	AA	1274	G	C5-C6-O6	-6.78	124.53	128.60
1	AA	1497	G	C5-C6-O6	-6.78	124.53	128.60
53	B5	320	G	N1-C6-O6	6.78	123.97	119.90
53	B5	814	U	O4'-C1'-N1	6.78	113.62	108.20
53	B5	2105	G	N1-C6-O6	6.78	123.97	119.90
53	B5	2939	G	O4'-C1'-N9	6.78	113.62	108.20
53	B5	3339	A	O4'-C1'-N9	6.78	113.62	108.20
1	AA	588	U	O4'-C1'-N1	6.78	113.62	108.20
1	AA	635	A	O4'-C1'-N9	6.78	113.62	108.20
1	AA	1518	U	C4'-C3'-C2'	6.78	109.38	102.60
1	AA	1522	A	P-O3'-C3'	-6.78	111.57	119.70
1	AA	1547	C	C1'-C2'-O2'	6.78	130.92	110.60
1	AA	1681	C	O4'-C1'-N1	6.78	113.62	108.20
53	B5	141	C	O4'-C1'-N1	6.78	113.62	108.20
53	B5	3354	U	O4'-C1'-N1	6.78	113.62	108.20
1	AA	1484	G	N1-C6-O6	6.77	123.97	119.90
19	A7	3	G	C8-N9-C4	-6.77	103.69	106.40
19	A7	62	A	C5'-C4'-C3'	-6.77	105.17	116.00
53	B5	1528	G	N1-C6-O6	6.77	123.96	119.90
53	B5	1825	G	C5-C6-O6	-6.77	124.54	128.60
1	AA	1034	G	N1-C6-O6	6.77	123.96	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	2614	G	C5-C6-O6	-6.77	124.54	128.60
1	AA	893	U	O4'-C1'-N1	6.77	113.61	108.20
53	B5	1877	U	O4'-C1'-N1	6.77	113.61	108.20
1	AA	660	G	N1-C6-O6	6.77	123.96	119.90
1	AA	1194	G	N1-C6-O6	6.77	123.96	119.90
52	B4	75	G	N1-C6-O6	6.77	123.96	119.90
53	B5	78	U	O4'-C1'-N1	6.77	113.61	108.20
53	B5	1058	U	O4'-C1'-N1	6.77	113.61	108.20
53	B5	1184	A	O4'-C1'-N9	6.77	113.61	108.20
53	B5	1141	C	O4'-C1'-N1	6.77	113.61	108.20
53	B5	1421	G	C5-C6-O6	-6.77	124.54	128.60
53	B5	2617	U	O4'-C1'-N1	6.76	113.61	108.20
53	B5	36	C	O4'-C1'-N1	6.76	113.61	108.20
53	B5	977	U	O4'-C1'-N1	6.76	113.61	108.20
53	B5	1319	G	C5-C6-O6	-6.76	124.54	128.60
53	B5	1365	G	C5-C6-O6	-6.76	124.54	128.60
1	AA	1209	G	O4'-C1'-N9	6.76	113.61	108.20
53	B5	727	G	C5-C6-O6	-6.76	124.54	128.60
53	B5	900	G	N1-C6-O6	6.76	123.96	119.90
53	B5	1411	C	P-O3'-C3'	6.76	127.81	119.70
53	B5	1581	C	P-O3'-C3'	6.76	127.81	119.70
53	B5	2499	U	O4'-C1'-N1	6.76	113.61	108.20
1	AA	87	C	O4'-C1'-N1	6.76	113.61	108.20
51	B3	111	U	O4'-C1'-N1	6.76	113.61	108.20
53	B5	712	G	O4'-C1'-N9	6.76	113.61	108.20
53	B5	994	G	O4'-C1'-N9	6.76	113.61	108.20
1	AA	1456	G	N1-C6-O6	6.76	123.95	119.90
19	A7	75	C	C4-C5-C6	-6.76	114.02	117.40
53	B5	1123	U	O4'-C1'-N1	6.76	113.61	108.20
53	B5	2426	U	O4'-C1'-N1	6.76	113.61	108.20
1	AA	722	G	N1-C6-O6	6.76	123.95	119.90
1	AA	1285	G	N1-C6-O6	6.76	123.95	119.90
52	B4	150	G	C5-C6-O6	-6.76	124.55	128.60
53	B5	800	G	C5-C6-O6	-6.76	124.55	128.60
53	B5	2693	C	O4'-C1'-N1	6.76	113.61	108.20
52	B4	1	A	O4'-C1'-N9	6.75	113.60	108.20
1	AA	944	U	O4'-C1'-N1	6.75	113.60	108.20
53	B5	749	C	O4'-C1'-N1	6.75	113.60	108.20
53	B5	1531	C	O4'-C1'-N1	6.75	113.60	108.20
53	B5	2121	G	N1-C6-O6	6.75	123.95	119.90
53	B5	2850	G	N1-C6-O6	6.75	123.95	119.90
1	AA	70	C	O4'-C1'-N1	6.75	113.60	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	426	G	N1-C6-O6	6.75	123.95	119.90
1	AA	747	C	O4'-C1'-N1	6.75	113.60	108.20
1	AA	826	U	O4'-C1'-N1	6.75	113.60	108.20
1	AA	903	G	N1-C6-O6	6.75	123.95	119.90
1	AA	1552	U	C4'-C3'-C2'	6.75	109.35	102.60
53	B5	897	U	O4'-C1'-N1	6.75	113.60	108.20
53	B5	2437	G	N1-C6-O6	6.75	123.95	119.90
1	AA	29	U	O4'-C1'-N1	6.75	113.60	108.20
53	B5	3088	G	C5-C6-O6	-6.75	124.55	128.60
1	AA	461	G	N1-C6-O6	6.75	123.95	119.90
1	AA	1426	G	N1-C6-O6	6.75	123.95	119.90
1	AA	1561	C	O4'-C1'-N1	6.75	113.60	108.20
51	B3	16	U	O4'-C1'-N1	6.75	113.60	108.20
1	AA	1516	C	O4'-C1'-N1	6.75	113.60	108.20
1	AA	1517	U	O4'-C4'-C3'	-6.75	97.25	104.00
53	B5	2173	U	O4'-C1'-N1	6.75	113.60	108.20
1	AA	1198	A	O4'-C1'-N9	6.74	113.59	108.20
1	AA	1208	G	N1-C6-O6	6.74	123.94	119.90
1	AA	667	U	O4'-C1'-N1	6.74	113.59	108.20
1	AA	895	G	N1-C6-O6	6.74	123.94	119.90
53	B5	1164	G	C5-C6-O6	-6.74	124.56	128.60
1	AA	768	C	O4'-C1'-N1	6.74	113.59	108.20
53	B5	631	U	O4'-C1'-N1	6.74	113.59	108.20
53	B5	2912	G	N1-C6-O6	6.74	123.94	119.90
1	AA	179	A	O4'-C1'-N9	6.74	113.59	108.20
1	AA	540	G	N1-C6-O6	6.74	123.94	119.90
1	AA	546	U	O4'-C1'-N1	6.74	113.59	108.20
1	AA	1013	G	C5-C6-O6	-6.74	124.56	128.60
1	AA	1582	G	N1-C6-O6	6.74	123.94	119.90
53	B5	616	G	N1-C6-O6	6.74	123.94	119.90
53	B5	1668	G	C5-C6-O6	-6.74	124.56	128.60
53	B5	2724	U	O4'-C1'-N1	6.74	113.59	108.20
53	B5	3107	U	O4'-C1'-N1	6.74	113.59	108.20
53	B5	1429	G	C5-C6-O6	-6.73	124.56	128.60
53	B5	1544	G	C5-C6-O6	-6.73	124.56	128.60
53	B5	2418	G	N1-C6-O6	6.73	123.94	119.90
1	AA	551	G	N1-C6-O6	6.73	123.94	119.90
1	AA	1144	A	C5-C6-N6	-6.73	118.32	123.70
52	B4	121	U	O4'-C1'-N1	6.73	113.58	108.20
53	B5	872	U	O4'-C1'-N1	6.73	113.58	108.20
53	B5	2234	G	N1-C6-O6	6.73	123.94	119.90
53	B5	2465	G	C5-C6-O6	-6.73	124.56	128.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	334	G	N1-C6-O6	6.73	123.94	119.90
1	AA	1221	U	O4'-C1'-N1	6.73	113.58	108.20
53	B5	101	G	N1-C6-O6	6.73	123.94	119.90
53	B5	176	G	O4'-C1'-N9	6.73	113.58	108.20
53	B5	946	U	O4'-C1'-N1	6.73	113.58	108.20
1	AA	347	G	C5-C6-O6	-6.72	124.56	128.60
1	AA	1755	G	O4'-C1'-N9	6.72	113.58	108.20
19	A7	47	U	C2-N3-C4	-6.72	122.97	127.00
53	B5	1530	U	O4'-C1'-N1	6.72	113.58	108.20
53	B5	2962	U	O4'-C1'-N1	6.72	113.58	108.20
1	AA	42	G	N1-C6-O6	6.72	123.93	119.90
51	B3	63	G	O4'-C1'-N9	6.72	113.58	108.20
53	B5	1568	U	O4'-C1'-N1	6.72	113.58	108.20
53	B5	2307	G	C5-C6-O6	-6.72	124.57	128.60
53	B5	2570	U	O4'-C1'-N1	6.72	113.58	108.20
1	AA	154	G	O4'-C1'-N9	6.72	113.58	108.20
1	AA	95	G	N1-C6-O6	6.72	123.93	119.90
1	AA	1146	G	N1-C6-O6	6.72	123.93	119.90
52	B4	94	C	O4'-C1'-N1	6.72	113.58	108.20
53	B5	83	U	O4'-C1'-N1	6.72	113.58	108.20
53	B5	280	U	O4'-C1'-N1	6.72	113.58	108.20
53	B5	742	G	N1-C6-O6	6.72	123.93	119.90
53	B5	1916	U	O4'-C1'-N1	6.72	113.58	108.20
53	B5	2623	G	C5-C6-O6	-6.72	124.57	128.60
53	B5	2760	C	O4'-C1'-N1	6.72	113.58	108.20
53	B5	3074	G	O4'-C1'-N9	6.72	113.58	108.20
53	B5	2374	C	O4'-C1'-N1	6.72	113.57	108.20
1	AA	411	C	O4'-C1'-N1	6.72	113.57	108.20
53	B5	1214	U	O4'-C1'-N1	6.72	113.57	108.20
53	B5	2227	C	O4'-C1'-N1	6.72	113.57	108.20
1	AA	958	U	O4'-C1'-N1	6.71	113.57	108.20
1	AA	1329	C	O4'-C1'-N1	6.71	113.57	108.20
53	B5	1517	G	N1-C6-O6	6.71	123.93	119.90
1	AA	571	G	N1-C6-O6	6.71	123.93	119.90
1	AA	586	G	N1-C6-O6	6.71	123.93	119.90
1	AA	1010	G	N1-C6-O6	6.71	123.93	119.90
1	AA	1053	U	O4'-C1'-N1	6.71	113.57	108.20
53	B5	1564	U	O4'-C1'-N1	6.71	113.57	108.20
53	B5	2125	A	O4'-C1'-N9	6.71	113.57	108.20
53	B5	509	U	O4'-C1'-N1	6.71	113.57	108.20
53	B5	549	U	O4'-C1'-N1	6.71	113.57	108.20
53	B5	2565	U	O4'-C1'-N1	6.71	113.57	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	196	G	N1-C6-O6	6.71	123.92	119.90
1	AA	1745	G	N1-C6-O6	6.71	123.92	119.90
52	B4	156	U	O4'-C1'-N1	6.71	113.57	108.20
53	B5	237	G	C5-C6-O6	-6.71	124.58	128.60
53	B5	591	U	O4'-C1'-N1	6.71	113.57	108.20
53	B5	2153	U	O4'-C1'-N1	6.71	113.57	108.20
53	B5	2992	U	O4'-C1'-N1	6.71	113.57	108.20
1	AA	1592	G	N1-C6-O6	6.71	123.92	119.90
1	AA	1645	U	O4'-C1'-N1	6.71	113.56	108.20
53	B5	1066	G	N1-C6-O6	6.71	123.92	119.90
53	B5	2592	G	N1-C6-O6	6.71	123.92	119.90
1	AA	921	G	N1-C6-O6	6.71	123.92	119.90
53	B5	2253	G	N1-C6-O6	6.71	123.92	119.90
53	B5	2699	G	N1-C6-O6	6.71	123.92	119.90
53	B5	2815	G	N1-C6-O6	6.71	123.92	119.90
53	B5	2875	U	O4'-C1'-N1	6.71	113.56	108.20
1	AA	292	U	O4'-C1'-N1	6.70	113.56	108.20
1	AA	1559	U	O4'-C1'-N1	6.70	113.56	108.20
53	B5	2429	G	N1-C6-O6	6.70	123.92	119.90
51	B3	103	U	O4'-C1'-N1	6.70	113.56	108.20
1	AA	778	G	N1-C6-O6	6.70	123.92	119.90
1	AA	831	U	O4'-C1'-N1	6.70	113.56	108.20
1	AA	1546	G	C2'-C3'-O3'	-6.70	94.76	109.50
1	AA	1769	U	O4'-C1'-N1	6.70	113.56	108.20
1	AA	1795	A	O4'-C1'-N9	6.70	113.56	108.20
53	B5	2607	G	N1-C6-O6	6.70	123.92	119.90
53	B5	2620	G	C5-C6-O6	-6.70	124.58	128.60
1	AA	797	G	OP1-P-OP2	-6.70	109.55	119.60
1	AA	1549	U	C2'-C3'-O3'	6.70	124.42	113.70
19	A7	65	G	N7-C8-N9	6.70	116.45	113.10
53	B5	505	G	O4'-C1'-N9	6.70	113.56	108.20
53	B5	2901	G	N1-C6-O6	6.70	123.92	119.90
53	B5	2266	U	O4'-C1'-N1	6.70	113.56	108.20
1	AA	268	C	O4'-C1'-N1	6.70	113.56	108.20
19	A7	14	A	C5-N7-C8	6.70	107.25	103.90
30	BF	138	TYR	CB-CG-CD1	6.70	125.02	121.00
52	B4	50	C	O4'-C1'-N1	6.70	113.56	108.20
53	B5	494	G	O4'-C1'-N9	6.70	113.56	108.20
53	B5	714	G	C5-C6-O6	-6.70	124.58	128.60
53	B5	1463	U	O4'-C1'-N1	6.70	113.56	108.20
53	B5	3061	G	C5-C6-O6	-6.70	124.58	128.60
53	B5	564	G	N1-C6-O6	6.69	123.92	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1577	U	O4'-C1'-N1	6.69	113.55	108.20
24	B9	14	TYR	CB-CG-CD2	6.69	125.02	121.00
53	B5	2190	U	O4'-C1'-N1	6.69	113.55	108.20
1	AA	857	U	O5'-C5'-C4'	6.69	124.41	111.70
1	AA	1540	G	N1-C6-O6	6.69	123.91	119.90
52	B4	29	U	O4'-C1'-N1	6.69	113.55	108.20
53	B5	2442	G	N1-C6-O6	6.69	123.91	119.90
53	B5	2880	U	O4'-C1'-N1	6.69	113.55	108.20
53	B5	3199	G	N1-C6-O6	6.69	123.91	119.90
1	AA	1260	G	N1-C6-O6	6.69	123.91	119.90
53	B5	1737	U	O4'-C1'-N1	6.69	113.55	108.20
53	B5	147	U	O4'-C1'-N1	6.69	113.55	108.20
53	B5	1487	G	C5-C6-O6	-6.69	124.59	128.60
53	B5	3116	G	O4'-C1'-N9	6.69	113.55	108.20
1	AA	484	C	O4'-C1'-N1	6.68	113.55	108.20
1	AA	907	U	O4'-C1'-N1	6.68	113.55	108.20
1	AA	1464	G	N1-C6-O6	6.68	123.91	119.90
1	AA	1690	G	N1-C6-O6	6.68	123.91	119.90
19	A7	14	A	N7-C8-N9	-6.68	110.46	113.80
53	B5	3264	G	C5-C6-O6	-6.68	124.59	128.60
1	AA	1161	G	N1-C6-O6	6.68	123.91	119.90
1	AA	1382	G	O4'-C1'-N9	6.68	113.55	108.20
1	AA	1671	G	N1-C6-O6	6.68	123.91	119.90
19	A7	4	G	C6-C5-N7	6.68	134.41	130.40
42	BR	94	TYR	CB-CG-CD1	-6.68	116.99	121.00
53	B5	354	U	O4'-C1'-N1	6.68	113.55	108.20
53	B5	1380	G	O4'-C1'-N9	6.68	113.55	108.20
53	B5	2534	G	N1-C6-O6	6.68	123.91	119.90
1	AA	1344	U	O4'-C1'-N1	6.68	113.55	108.20
53	B5	2618	G	C5-C6-O6	-6.68	124.59	128.60
1	AA	1506	U	O4'-C1'-N1	6.68	113.54	108.20
1	AA	1636	G	C1'-O4'-C4'	-6.68	104.56	109.90
53	B5	2158	A	P-O3'-C3'	6.68	127.72	119.70
53	B5	2809	C	O4'-C1'-N1	6.68	113.54	108.20
53	B5	3115	C	O4'-C1'-N1	6.68	113.54	108.20
53	B5	1172	G	N1-C6-O6	6.68	123.91	119.90
53	B5	1819	U	O4'-C1'-N1	6.68	113.54	108.20
53	B5	2986	U	O4'-C1'-N1	6.68	113.54	108.20
53	B5	3019	U	O4'-C1'-N1	6.68	113.54	108.20
1	AA	106	U	O4'-C1'-N1	6.68	113.54	108.20
53	B5	3098	G	C5-C6-O6	-6.68	124.59	128.60
1	AA	1632	C	P-O3'-C3'	-6.67	111.69	119.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	263	C	O4'-C1'-N1	6.67	113.54	108.20
53	B5	1408	G	N1-C6-O6	6.67	123.91	119.90
53	B5	1470	U	O4'-C1'-N1	6.67	113.54	108.20
53	B5	3284	G	C5-C6-O6	-6.67	124.60	128.60
53	B5	1612	A	O4'-C1'-N9	6.67	113.54	108.20
1	AA	1249	U	O4'-C1'-N1	6.67	113.54	108.20
1	AA	1448	U	O4'-C1'-N1	6.67	113.54	108.20
53	B5	410	U	O4'-C1'-N1	6.67	113.54	108.20
53	B5	996	U	O4'-C1'-N1	6.67	113.54	108.20
1	AA	496	G	N1-C6-O6	6.67	123.90	119.90
53	B5	155	G	C5-C6-O6	-6.67	124.60	128.60
53	B5	739	G	O4'-C1'-N9	6.67	113.54	108.20
53	B5	2469	G	C5-C6-O6	-6.67	124.60	128.60
51	B3	5	G	O4'-C1'-N9	6.67	113.53	108.20
53	B5	471	U	O4'-C1'-N1	6.67	113.53	108.20
53	B5	524	C	O4'-C1'-N1	6.67	113.53	108.20
53	B5	619	A	O4'-C1'-N9	6.67	113.53	108.20
53	B5	625	G	N1-C6-O6	6.67	123.90	119.90
53	B5	627	U	O4'-C1'-N1	6.67	113.53	108.20
53	B5	609	G	C5-C6-O6	-6.67	124.60	128.60
53	B5	2610	G	N1-C6-O6	6.67	123.90	119.90
53	B5	2806	U	O4'-C1'-N1	6.67	113.53	108.20
53	B5	733	G	N1-C6-O6	6.67	123.90	119.90
53	B5	2599	U	O4'-C1'-N1	6.67	113.53	108.20
1	AA	633	U	O4'-C1'-N1	6.66	113.53	108.20
1	AA	834	G	O4'-C1'-N9	6.66	113.53	108.20
51	B3	46	A	O4'-C1'-N9	6.66	113.53	108.20
53	B5	632	G	N1-C6-O6	6.66	123.90	119.90
1	AA	699	U	O4'-C1'-N1	6.66	113.53	108.20
1	AA	610	G	N1-C6-O6	6.66	123.90	119.90
16	AS	97	TYR	CB-CG-CD2	-6.66	117.00	121.00
53	B5	860	G	N1-C6-O6	6.66	123.90	119.90
53	B5	1651	U	O4'-C1'-N1	6.66	113.53	108.20
1	AA	1200	A	C6-C5-N7	6.66	136.96	132.30
52	B4	31	G	O4'-C1'-N9	6.66	113.53	108.20
53	B5	1683	A	O4'-C1'-N9	6.66	113.53	108.20
53	B5	1874	A	O4'-C1'-N9	6.66	113.53	108.20
53	B5	1942	U	O4'-C1'-N1	6.66	113.53	108.20
1	AA	1233	G	N1-C6-O6	6.66	123.89	119.90
53	B5	618	C	O4'-C1'-N1	6.66	113.53	108.20
1	AA	702	G	N1-C6-O6	6.66	123.89	119.90
19	A7	18	G	C6-C5-N7	6.66	134.39	130.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
51	B3	118	U	O4'-C1'-N1	6.66	113.52	108.20
53	B5	3045	G	N1-C6-O6	6.66	123.89	119.90
53	B5	3310	A	C5-C6-N6	-6.66	118.38	123.70
53	B5	2886	U	O4'-C1'-N1	6.65	113.52	108.20
53	B5	2915	U	O4'-C1'-N1	6.65	113.52	108.20
53	B5	3075	G	O4'-C1'-N9	6.65	113.52	108.20
1	AA	1723	U	O4'-C1'-N1	6.65	113.52	108.20
53	B5	58	G	O4'-C1'-N9	6.65	113.52	108.20
53	B5	2791	G	O4'-C1'-N9	6.65	113.52	108.20
53	B5	2839	G	N1-C6-O6	6.65	123.89	119.90
53	B5	3346	U	O4'-C1'-N1	6.65	113.52	108.20
51	B3	85	G	N1-C6-O6	6.65	123.89	119.90
53	B5	2945	G	C5-C6-O6	-6.65	124.61	128.60
1	AA	1750	U	O4'-C1'-N1	6.65	113.52	108.20
53	B5	137	G	O4'-C1'-N9	6.65	113.52	108.20
1	AA	830	U	O4'-C1'-N1	6.64	113.52	108.20
52	B4	153	U	O4'-C1'-N1	6.64	113.52	108.20
53	B5	2324	A	O4'-C1'-N9	6.64	113.52	108.20
53	B5	3286	G	N1-C6-O6	6.64	123.89	119.90
1	AA	23	G	N1-C6-O6	6.64	123.89	119.90
1	AA	822	U	O4'-C1'-N1	6.64	113.51	108.20
1	AA	886	U	O4'-C1'-N1	6.64	113.51	108.20
19	A7	15	G	C1'-O4'-C4'	-6.64	104.59	109.90
53	B5	189	G	O4'-C1'-N9	6.64	113.51	108.20
1	AA	1610	U	P-O3'-C3'	6.64	127.67	119.70
53	B5	1469	C	O4'-C1'-N1	6.64	113.51	108.20
53	B5	1570	U	O4'-C1'-N1	6.64	113.51	108.20
53	B5	1914	G	N1-C6-O6	6.64	123.88	119.90
53	B5	2300	G	N1-C6-O6	6.64	123.88	119.90
1	AA	73	U	O4'-C1'-N1	6.64	113.51	108.20
53	B5	301	G	O4'-C1'-N9	6.64	113.51	108.20
1	AA	404	G	N1-C6-O6	6.64	123.88	119.90
1	AA	554	C	O4'-C1'-N1	6.64	113.51	108.20
1	AA	48	G	O4'-C1'-N9	6.63	113.51	108.20
1	AA	1092	U	O4'-C1'-N1	6.63	113.51	108.20
18	AT	18	TYR	CB-CG-CD2	-6.63	117.02	121.00
53	B5	1379	G	C5-C6-O6	-6.63	124.62	128.60
53	B5	1387	G	C5-C6-O6	-6.63	124.62	128.60
53	B5	2470	C	C2-N1-C1'	6.63	126.10	118.80
53	B5	2883	U	O4'-C1'-N1	6.63	113.51	108.20
1	AA	289	U	O4'-C1'-N1	6.63	113.51	108.20
53	B5	70	A	P-O3'-C3'	6.63	127.66	119.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	91	G	N1-C6-O6	6.63	123.88	119.90
1	AA	1000	U	O4'-C1'-N1	6.63	113.50	108.20
1	AA	1150	G	N1-C6-O6	6.63	123.88	119.90
1	AA	1315	G	N1-C6-O6	6.63	123.88	119.90
1	AA	1706	U	O4'-C1'-N1	6.63	113.50	108.20
53	B5	328	U	O4'-C1'-N1	6.63	113.50	108.20
53	B5	632	G	O4'-C1'-N9	6.63	113.50	108.20
53	B5	1409	G	N1-C6-O6	6.63	123.88	119.90
53	B5	2343	C	O4'-C1'-N1	6.63	113.50	108.20
53	B5	2575	G	O4'-C1'-N9	6.63	113.51	108.20
1	AA	858	G	N7-C8-N9	6.63	116.42	113.10
53	B5	1811	G	O4'-C1'-N9	6.63	113.50	108.20
53	B5	2497	U	O4'-C1'-N1	6.63	113.50	108.20
53	B5	3149	G	C5-C6-O6	-6.63	124.62	128.60
1	AA	976	A	O4'-C1'-N9	6.63	113.50	108.20
1	AA	1224	G	N1-C6-O6	6.63	123.88	119.90
53	B5	795	G	C5-C6-O6	-6.63	124.62	128.60
53	B5	1122	U	O4'-C1'-N1	6.63	113.50	108.20
53	B5	1217	A	O4'-C1'-N9	6.63	113.50	108.20
53	B5	2427	U	O4'-C1'-N1	6.63	113.50	108.20
53	B5	2505	U	O4'-C1'-N1	6.63	113.50	108.20
1	AA	467	G	C5-C6-O6	-6.63	124.62	128.60
1	AA	1599	G	N1-C6-O6	6.63	123.88	119.90
53	B5	563	U	O4'-C1'-N1	6.63	113.50	108.20
53	B5	1339	C	O4'-C1'-N1	6.63	113.50	108.20
53	B5	1514	G	N1-C6-O6	6.62	123.88	119.90
52	B4	14	C	O4'-C1'-N1	6.62	113.50	108.20
53	B5	22	G	C5-C6-O6	-6.62	124.63	128.60
53	B5	1295	G	N1-C6-O6	6.62	123.87	119.90
53	B5	2559	A	P-O3'-C3'	6.62	127.65	119.70
1	AA	83	G	N1-C6-O6	6.62	123.87	119.90
1	AA	418	G	N1-C6-O6	6.62	123.87	119.90
1	AA	1549	U	C1'-C2'-O2'	6.62	130.46	110.60
52	B4	78	G	C5-C6-O6	-6.62	124.63	128.60
53	B5	286	U	O4'-C1'-N1	6.62	113.50	108.20
53	B5	601	U	P-O3'-C3'	6.62	127.64	119.70
53	B5	2252	A	O4'-C1'-N9	6.62	113.50	108.20
1	AA	1263	G	N1-C6-O6	6.62	123.87	119.90
51	B3	84	G	O4'-C1'-N9	6.62	113.50	108.20
53	B5	623	U	O4'-C1'-N1	6.62	113.49	108.20
53	B5	939	U	O4'-C1'-N1	6.62	113.50	108.20
53	B5	1400	G	N1-C6-O6	6.62	123.87	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	1735	G	C5-C6-O6	-6.62	124.63	128.60
53	B5	1735	G	O4'-C1'-N9	6.62	113.50	108.20
53	B5	3359	A	O4'-C1'-N9	6.62	113.49	108.20
53	B5	3394	U	O4'-C1'-N1	6.62	113.50	108.20
1	AA	1572	G	N1-C6-O6	6.62	123.87	119.90
53	B5	1282	G	C5-C6-O6	-6.62	124.63	128.60
1	AA	1563	C	P-O3'-C3'	6.62	127.64	119.70
53	B5	74	G	O4'-C1'-N9	6.62	113.49	108.20
53	B5	620	U	O4'-C1'-N1	6.62	113.49	108.20
53	B5	212	G	O4'-C1'-N9	6.61	113.49	108.20
53	B5	1062	A	O4'-C1'-N9	6.61	113.49	108.20
53	B5	3278	U	O4'-C1'-N1	6.61	113.49	108.20
53	B5	1586	G	C5-C6-O6	-6.61	124.63	128.60
1	AA	671	G	N1-C6-O6	6.61	123.87	119.90
53	B5	857	G	N1-C6-O6	6.61	123.87	119.90
53	B5	1653	G	C5-C6-O6	-6.61	124.63	128.60
53	B5	2184	U	O4'-C1'-N1	6.61	113.49	108.20
19	A7	4	G	C4-C5-C6	-6.61	114.83	118.80
53	B5	218	G	C5-C6-O6	-6.61	124.64	128.60
53	B5	535	G	N1-C6-O6	6.61	123.86	119.90
53	B5	1485	G	N1-C6-O6	6.61	123.87	119.90
53	B5	2807	U	O4'-C1'-N1	6.61	113.49	108.20
1	AA	462	G	N1-C6-O6	6.61	123.86	119.90
53	B5	44	U	O4'-C1'-N1	6.61	113.48	108.20
53	B5	257	U	O4'-C1'-N1	6.61	113.48	108.20
53	B5	467	U	O4'-C1'-N1	6.61	113.48	108.20
53	B5	953	G	C5-C6-O6	-6.61	124.64	128.60
53	B5	1600	U	O4'-C1'-N1	6.61	113.48	108.20
53	B5	1880	U	O4'-C1'-N1	6.61	113.48	108.20
53	B5	2650	U	O4'-C1'-N1	6.61	113.48	108.20
1	AA	1182	U	O4'-C1'-N1	6.60	113.48	108.20
53	B5	258	G	C5-C6-O6	-6.60	124.64	128.60
53	B5	2391	G	N1-C6-O6	6.60	123.86	119.90
53	B5	3326	G	O4'-C1'-N9	6.60	113.48	108.20
1	AA	739	G	N1-C6-O6	6.60	123.86	119.90
53	B5	2591	A	O4'-C1'-N9	6.60	113.48	108.20
1	AA	12	U	O4'-C1'-N1	6.60	113.48	108.20
1	AA	24	U	O4'-C1'-N1	6.60	113.48	108.20
52	B4	58	G	N1-C6-O6	6.60	123.86	119.90
53	B5	940	G	N1-C6-O6	6.60	123.86	119.90
53	B5	3159	C	O4'-C1'-N1	6.60	113.48	108.20
1	AA	857	U	C5'-C4'-C3'	6.60	126.56	116.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
19	A7	36	A	C4'-C3'-C2'	-6.60	96.00	102.60
53	B5	1845	G	C5-C6-O6	-6.60	124.64	128.60
1	AA	1178	U	P-O3'-C3'	6.60	127.62	119.70
53	B5	495	G	P-O3'-C3'	6.60	127.62	119.70
53	B5	1876	U	O4'-C1'-N1	6.60	113.48	108.20
53	B5	2537	U	O4'-C1'-N1	6.60	113.48	108.20
1	AA	357	G	O4'-C1'-N9	6.60	113.48	108.20
1	AA	1721	U	O4'-C1'-N1	6.60	113.48	108.20
53	B5	2522	G	N1-C6-O6	6.60	123.86	119.90
1	AA	496	G	O4'-C1'-N9	6.59	113.48	108.20
1	AA	971	G	N1-C6-O6	6.59	123.86	119.90
1	AA	1597	C	O4'-C1'-N1	6.59	113.48	108.20
52	B4	25	G	C5-C6-O6	-6.59	124.64	128.60
53	B5	429	U	O4'-C1'-N1	6.59	113.47	108.20
53	B5	2327	U	O4'-C1'-N1	6.59	113.48	108.20
53	B5	2874	G	N1-C6-O6	6.59	123.86	119.90
53	B5	1927	G	C5-C6-O6	-6.59	124.64	128.60
1	AA	21	U	O4'-C1'-N1	6.59	113.47	108.20
1	AA	874	C	O4'-C1'-N1	6.59	113.47	108.20
1	AA	1601	U	O4'-C1'-N1	6.59	113.47	108.20
1	AA	945	U	O4'-C1'-N1	6.59	113.47	108.20
1	AA	1256	U	O4'-C1'-N1	6.59	113.47	108.20
1	AA	1697	G	N1-C6-O6	6.59	123.85	119.90
16	AS	97	TYR	CB-CG-CD1	6.59	124.95	121.00
53	B5	456	U	O4'-C1'-N1	6.59	113.47	108.20
53	B5	1261	G	O4'-C1'-N9	6.59	113.47	108.20
53	B5	1716	U	O4'-C1'-N1	6.59	113.47	108.20
53	B5	1861	G	C5-C6-O6	-6.59	124.65	128.60
1	AA	1391	G	N1-C6-O6	6.59	123.85	119.90
1	AA	1558	U	O4'-C1'-N1	6.59	113.47	108.20
53	B5	487	U	O4'-C1'-N1	6.59	113.47	108.20
53	B5	520	U	O4'-C1'-N1	6.59	113.47	108.20
53	B5	3036	G	O4'-C1'-N9	6.59	113.47	108.20
1	AA	430	G	N1-C6-O6	6.58	123.85	119.90
53	B5	458	U	O4'-C1'-N1	6.58	113.47	108.20
53	B5	1139	G	C5-C6-O6	-6.58	124.65	128.60
53	B5	2586	G	N1-C6-O6	6.58	123.85	119.90
1	AA	703	G	C5-C6-O6	-6.58	124.65	128.60
1	AA	1725	G	N1-C6-O6	6.58	123.85	119.90
53	B5	517	G	N1-C6-O6	6.58	123.85	119.90
53	B5	1565	G	C5-C6-O6	-6.58	124.65	128.60
53	B5	2997	G	C5-C6-O6	-6.58	124.65	128.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	508	U	O4'-C1'-N1	6.58	113.47	108.20
49	BY	47	TYR	CB-CG-CD1	-6.58	117.05	121.00
53	B5	459	G	N1-C6-O6	6.58	123.85	119.90
53	B5	942	U	O4'-C1'-N1	6.58	113.46	108.20
1	AA	984	G	N1-C6-O6	6.58	123.85	119.90
53	B5	80	G	N1-C6-O6	6.58	123.85	119.90
53	B5	404	G	C5-C6-O6	-6.58	124.65	128.60
53	B5	2907	G	N1-C6-O6	6.58	123.85	119.90
1	AA	457	G	N1-C6-O6	6.58	123.84	119.90
53	B5	859	G	N1-C6-O6	6.58	123.84	119.90
53	B5	1158	A	O4'-C1'-N9	6.58	113.46	108.20
53	B5	2731	U	O4'-C1'-N1	6.58	113.46	108.20
53	B5	3387	U	O4'-C1'-N1	6.58	113.46	108.20
1	AA	296	U	O4'-C1'-N1	6.57	113.46	108.20
1	AA	1603	G	N1-C6-O6	6.57	123.84	119.90
53	B5	347	G	N1-C6-O6	6.57	123.84	119.90
53	B5	421	G	C5-C6-O6	-6.57	124.66	128.60
53	B5	2261	G	N1-C6-O6	6.57	123.84	119.90
1	AA	195	G	N1-C6-O6	6.57	123.84	119.90
1	AA	1409	G	C5-C6-O6	-6.57	124.66	128.60
1	AA	1715	G	N1-C6-O6	6.57	123.84	119.90
1	AA	828	U	O4'-C1'-N1	6.57	113.46	108.20
19	A7	12	U	O4'-C4'-C3'	6.57	111.36	106.10
19	A7	28	C	N3-C4-C5	6.57	124.53	121.90
53	B5	76	G	C5-C6-O6	-6.57	124.66	128.60
53	B5	139	G	C5-C6-O6	-6.57	124.66	128.60
53	B5	1269	U	P-O3'-C3'	6.57	127.59	119.70
53	B5	2481	G	O4'-C1'-N9	6.57	113.46	108.20
53	B5	2909	U	O4'-C1'-N1	6.57	113.46	108.20
44	BT	60	TYR	CB-CG-CD2	-6.57	117.06	121.00
52	B4	148	G	O4'-C1'-N9	6.57	113.45	108.20
53	B5	1692	U	O4'-C1'-N1	6.57	113.45	108.20
1	AA	205	U	O4'-C1'-N1	6.57	113.45	108.20
1	AA	1727	C	O4'-C1'-N1	6.57	113.45	108.20
53	B5	2154	U	O4'-C1'-N1	6.57	113.45	108.20
53	B5	2380	U	O4'-C1'-N1	6.57	113.45	108.20
53	B5	3040	A	O4'-C1'-N9	6.57	113.45	108.20
53	B5	1468	A	O4'-C1'-N9	6.56	113.45	108.20
1	AA	248	U	O4'-C1'-N1	6.56	113.45	108.20
1	AA	1713	G	N1-C6-O6	6.56	123.84	119.90
53	B5	1361	U	O4'-C1'-N1	6.56	113.45	108.20
53	B5	1505	C	O4'-C1'-N1	6.56	113.45	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1151	G	N1-C6-O6	6.56	123.84	119.90
1	AA	1402	G	N1-C6-O6	6.56	123.84	119.90
53	B5	421	G	P-O3'-C3'	6.56	127.57	119.70
53	B5	844	G	N1-C6-O6	6.56	123.84	119.90
53	B5	2379	U	O4'-C1'-N1	6.56	113.45	108.20
1	AA	727	U	O4'-C1'-N1	6.56	113.45	108.20
53	B5	1743	G	N1-C6-O6	6.56	123.84	119.90
1	AA	1048	G	O4'-C1'-N9	6.56	113.45	108.20
1	AA	1688	G	N1-C6-O6	6.56	123.83	119.90
1	AA	1763	A	O4'-C1'-N9	6.56	113.45	108.20
53	B5	2123	G	O4'-C1'-N9	6.56	113.45	108.20
53	B5	3247	G	N1-C6-O6	6.56	123.83	119.90
1	AA	330	G	N1-C6-O6	6.56	123.83	119.90
1	AA	1152	G	C5-C6-O6	-6.56	124.67	128.60
52	B4	102	U	O4'-C1'-N1	6.56	113.44	108.20
53	B5	167	U	O4'-C1'-N1	6.56	113.44	108.20
53	B5	592	G	C5-C6-O6	-6.56	124.67	128.60
1	AA	866	G	OP1-P-OP2	6.55	129.43	119.60
1	AA	1520	U	C5'-C4'-O4'	6.55	116.97	109.10
53	B5	1464	G	C5-C6-O6	-6.55	124.67	128.60
53	B5	2767	U	O4'-C1'-N1	6.55	113.44	108.20
53	B5	2854	U	O4'-C1'-N1	6.55	113.44	108.20
1	AA	561	G	N1-C6-O6	6.55	123.83	119.90
1	AA	959	U	O4'-C1'-N1	6.55	113.44	108.20
1	AA	1137	G	N1-C6-O6	6.55	123.83	119.90
1	AA	1482	G	N1-C6-O6	6.55	123.83	119.90
48	BX	7	TYR	CB-CG-CD1	6.55	124.93	121.00
53	B5	425	G	C5-C6-O6	-6.55	124.67	128.60
53	B5	110	G	N1-C6-O6	6.55	123.83	119.90
53	B5	304	G	N1-C6-O6	6.55	123.83	119.90
53	B5	581	U	O4'-C1'-N1	6.55	113.44	108.20
53	B5	1264	G	N1-C6-O6	6.55	123.83	119.90
53	B5	1466	G	C5-C6-O6	-6.55	124.67	128.60
53	B5	2575	G	P-O3'-C3'	6.55	127.56	119.70
53	B5	3200	G	N1-C6-O6	6.55	123.83	119.90
53	B5	169	U	O4'-C1'-N1	6.55	113.44	108.20
53	B5	1533	U	O4'-C1'-N1	6.55	113.44	108.20
53	B5	2696	A	O4'-C1'-N9	6.55	113.44	108.20
1	AA	975	G	N1-C6-O6	6.55	123.83	119.90
1	AA	1105	G	C5-C6-O6	-6.55	124.67	128.60
1	AA	1175	G	N1-C6-O6	6.55	123.83	119.90
1	AA	1553	A	P-O5'-C5'	-6.55	110.43	120.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
19	A7	57	G	C5-N7-C8	-6.55	101.03	104.30
51	B3	30	G	C5-C6-O6	-6.55	124.67	128.60
53	B5	2816	G	N1-C6-O6	6.55	123.83	119.90
53	B5	3302	U	O4'-C1'-N1	6.55	113.44	108.20
1	AA	187	G	N1-C6-O6	6.54	123.83	119.90
53	B5	214	G	N1-C6-O6	6.54	123.83	119.90
1	AA	192	U	O4'-C1'-N1	6.54	113.44	108.20
1	AA	510	G	N1-C6-O6	6.54	123.83	119.90
1	AA	1486	G	C5-C6-O6	-6.54	124.67	128.60
53	B5	9	U	O4'-C1'-N1	6.54	113.43	108.20
53	B5	1641	U	O4'-C1'-N1	6.54	113.44	108.20
53	B5	2244	A	O4'-C1'-N9	6.54	113.44	108.20
53	B5	2482	U	O4'-C1'-N1	6.54	113.44	108.20
53	B5	528	U	O4'-C1'-N1	6.54	113.43	108.20
53	B5	1243	G	C5-C6-O6	-6.54	124.67	128.60
53	B5	1285	G	C5-C6-O6	-6.54	124.67	128.60
53	B5	2576	G	N1-C6-O6	6.54	123.83	119.90
53	B5	3344	A	O4'-C1'-N9	6.54	113.43	108.20
1	AA	920	U	O4'-C1'-N1	6.54	113.43	108.20
53	B5	783	A	O4'-C1'-N9	6.54	113.43	108.20
53	B5	1022	U	O4'-C1'-N1	6.54	113.43	108.20
1	AA	20	G	N1-C6-O6	6.54	123.82	119.90
1	AA	1271	A	O4'-C1'-N9	6.54	113.43	108.20
53	B5	651	G	N1-C6-O6	6.54	123.82	119.90
53	B5	1104	G	N1-C6-O6	6.54	123.82	119.90
53	B5	2138	A	P-O3'-C3'	6.54	127.54	119.70
1	AA	864	U	O4'-C1'-C2'	6.54	113.48	107.60
52	B4	6	U	O4'-C1'-N1	6.54	113.43	108.20
53	B5	935	U	O4'-C1'-N1	6.54	113.43	108.20
53	B5	2690	G	O4'-C1'-N9	6.54	113.43	108.20
1	AA	562	G	N1-C6-O6	6.53	123.82	119.90
53	B5	3272	U	O4'-C1'-N1	6.53	113.43	108.20
1	AA	1056	U	O4'-C1'-N1	6.53	113.43	108.20
53	B5	228	U	O4'-C1'-N1	6.53	113.43	108.20
53	B5	790	U	O4'-C1'-N1	6.53	113.42	108.20
1	AA	539	G	N1-C6-O6	6.53	123.82	119.90
1	AA	1586	G	O4'-C1'-N9	6.53	113.42	108.20
52	B4	87	G	C5-C6-O6	-6.53	124.68	128.60
53	B5	3046	A	O4'-C1'-N9	6.53	113.42	108.20
1	AA	274	G	N1-C6-O6	6.53	123.82	119.90
1	AA	1358	U	O4'-C1'-N1	6.53	113.42	108.20
51	B3	38	U	O4'-C1'-N1	6.53	113.42	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	531	G	O4'-C1'-N9	6.53	113.42	108.20
53	B5	777	U	O4'-C1'-N1	6.53	113.42	108.20
53	B5	1294	A	O4'-C1'-N9	6.53	113.42	108.20
1	AA	858	G	C8-N9-C4	-6.53	103.79	106.40
1	AA	941	G	O4'-C1'-N9	6.53	113.42	108.20
1	AA	1294	G	N1-C6-O6	6.53	123.81	119.90
53	B5	103	G	N1-C6-O6	6.53	123.82	119.90
53	B5	206	G	C5-C6-O6	-6.53	124.68	128.60
1	AA	1164	G	N1-C6-O6	6.52	123.81	119.90
53	B5	1389	G	C5-C6-O6	-6.52	124.69	128.60
53	B5	3181	G	N1-C6-O6	6.52	123.81	119.90
1	AA	1289	G	N1-C6-O6	6.52	123.81	119.90
1	AA	1380	G	N1-C6-O6	6.52	123.81	119.90
53	B5	532	A	O4'-C1'-N9	6.52	113.42	108.20
53	B5	1098	A	O4'-C1'-N9	6.52	113.42	108.20
53	B5	1512	U	O4'-C1'-N1	6.52	113.42	108.20
53	B5	1572	U	O4'-C1'-N1	6.52	113.42	108.20
53	B5	2947	G	N1-C6-O6	6.52	123.81	119.90
53	B5	3277	C	O4'-C1'-N1	6.52	113.42	108.20
53	B5	1248	C	O4'-C1'-N1	6.52	113.42	108.20
53	B5	2681	U	O4'-C1'-N1	6.52	113.42	108.20
1	AA	947	G	N1-C6-O6	6.52	123.81	119.90
1	AA	1346	G	O4'-C1'-N9	6.52	113.42	108.20
53	B5	451	U	O4'-C1'-N1	6.52	113.42	108.20
53	B5	1173	U	O4'-C1'-N1	6.52	113.42	108.20
53	B5	2410	U	O4'-C1'-N1	6.52	113.42	108.20
53	B5	2629	U	O4'-C1'-N1	6.52	113.42	108.20
53	B5	2701	U	O4'-C1'-N1	6.52	113.42	108.20
53	B5	2849	C	P-O3'-C3'	6.52	127.52	119.70
53	B5	173	G	O4'-C1'-N9	6.52	113.41	108.20
53	B5	262	U	O4'-C1'-N1	6.52	113.41	108.20
53	B5	275	U	O4'-C1'-N1	6.52	113.41	108.20
53	B5	716	G	O4'-C1'-N9	6.52	113.41	108.20
53	B5	2421	U	O4'-C1'-N1	6.52	113.41	108.20
53	B5	2773	C	O4'-C1'-N1	6.52	113.41	108.20
1	AA	864	U	O4'-C1'-N1	6.52	113.41	108.20
1	AA	873	U	O4'-C1'-N1	6.52	113.41	108.20
1	AA	1101	U	O4'-C1'-N1	6.52	113.41	108.20
53	B5	849	C	C6-N1-C1'	-6.52	112.98	120.80
1	AA	979	G	N1-C6-O6	6.51	123.81	119.90
19	A7	75	C	P-O3'-C3'	6.51	127.52	119.70
53	B5	792	G	N1-C6-O6	6.51	123.81	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	899	U	O4'-C1'-N1	6.51	113.41	108.20
53	B5	3014	U	O4'-C1'-N1	6.51	113.41	108.20
53	B5	3356	G	O4'-C1'-N9	6.51	113.41	108.20
53	B5	2517	U	O4'-C1'-N1	6.51	113.41	108.20
53	B5	3252	G	N1-C6-O6	6.51	123.81	119.90
1	AA	712	G	N1-C6-O6	6.51	123.81	119.90
19	A7	24	G	N3-C4-N9	6.51	129.91	126.00
53	B5	1520	G	O4'-C1'-N9	6.51	113.41	108.20
51	B3	74	A	O4'-C1'-N9	6.51	113.41	108.20
53	B5	921	A	O4'-C1'-N9	6.51	113.41	108.20
53	B5	2819	A	O4'-C1'-N9	6.51	113.41	108.20
53	B5	2855	U	O4'-C1'-N1	6.51	113.41	108.20
53	B5	3030	G	C5-C6-O6	-6.51	124.69	128.60
53	B5	2486	A	P-O3'-C3'	6.51	127.51	119.70
19	A7	5	A	O4'-C4'-C3'	6.51	111.31	106.10
19	A7	66	A	C5-N7-C8	6.51	107.15	103.90
52	B4	93	U	O4'-C1'-N1	6.51	113.41	108.20
52	B4	146	U	O4'-C1'-N1	6.51	113.40	108.20
53	B5	1636	U	O4'-C1'-N1	6.51	113.41	108.20
53	B5	1947	G	C5-C6-O6	-6.51	124.70	128.60
53	B5	2735	U	O4'-C1'-N1	6.51	113.41	108.20
53	B5	3197	G	C5-C6-O6	-6.51	124.70	128.60
51	B3	13	A	P-O3'-C3'	6.50	127.51	119.70
53	B5	370	U	O4'-C1'-N1	6.50	113.40	108.20
53	B5	510	G	N1-C6-O6	6.50	123.80	119.90
1	AA	835	U	O4'-C1'-N1	6.50	113.40	108.20
1	AA	478	A	O4'-C1'-N9	6.50	113.40	108.20
1	AA	672	U	O4'-C1'-N1	6.50	113.40	108.20
19	A7	35	A	C8-N9-C4	-6.50	103.20	105.80
53	B5	173	G	N1-C6-O6	6.50	123.80	119.90
53	B5	531	G	N1-C6-O6	6.50	123.80	119.90
53	B5	930	U	O4'-C1'-N1	6.50	113.40	108.20
53	B5	2187	G	N1-C6-O6	6.50	123.80	119.90
53	B5	2770	G	C5-C6-O6	-6.50	124.70	128.60
53	B5	3188	G	N1-C6-O6	6.50	123.80	119.90
1	AA	783	G	N1-C6-O6	6.50	123.80	119.90
1	AA	1111	G	N1-C6-O6	6.50	123.80	119.90
1	AA	1643	G	N1-C6-O6	6.50	123.80	119.90
1	AA	805	U	O4'-C1'-N1	6.50	113.40	108.20
1	AA	1378	U	O4'-C1'-N1	6.50	113.40	108.20
19	A7	51	G	N9-C1'-C2'	-6.50	104.85	112.00
53	B5	878	G	N1-C6-O6	6.50	123.80	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	1623	G	O4'-C1'-N9	6.50	113.40	108.20
53	B5	2470	C	O4'-C1'-N1	6.50	113.40	108.20
1	AA	234	G	N1-C6-O6	6.50	123.80	119.90
1	AA	1072	U	O4'-C1'-N1	6.50	113.40	108.20
1	AA	1313	G	N1-C6-O6	6.50	123.80	119.90
51	B3	79	U	O4'-C1'-N1	6.50	113.40	108.20
53	B5	947	G	N1-C6-O6	6.50	123.80	119.90
53	B5	1851	G	N1-C6-O6	6.50	123.80	119.90
1	AA	778	G	O4'-C1'-N9	6.49	113.39	108.20
1	AA	924	G	N1-C6-O6	6.49	123.80	119.90
53	B5	831	G	O4'-C1'-N9	6.49	113.39	108.20
53	B5	2945	G	O4'-C1'-N9	6.49	113.39	108.20
53	B5	3288	G	C5-C6-O6	-6.49	124.70	128.60
1	AA	743	U	O4'-C1'-N1	6.49	113.39	108.20
53	B5	2627	C	O4'-C1'-N1	6.49	113.39	108.20
1	AA	690	G	N1-C6-O6	6.49	123.79	119.90
1	AA	1399	G	N1-C6-O6	6.49	123.79	119.90
1	AA	1779	A	O4'-C1'-N9	6.49	113.39	108.20
53	B5	333	G	N1-C6-O6	6.49	123.79	119.90
53	B5	1719	G	C5-C6-O6	-6.49	124.71	128.60
53	B5	1860	G	N1-C6-O6	6.49	123.79	119.90
1	AA	211	U	O4'-C1'-N1	6.49	113.39	108.20
1	AA	815	G	N1-C6-O6	6.49	123.79	119.90
1	AA	1718	G	N1-C6-O6	6.49	123.79	119.90
1	AA	1735	G	N1-C6-O6	6.49	123.79	119.90
53	B5	3289	G	C5-C6-O6	-6.49	124.71	128.60
53	B5	59	G	C5-C6-O6	-6.49	124.71	128.60
53	B5	712	G	C5-C6-O6	-6.49	124.71	128.60
53	B5	1924	U	O4'-C1'-N1	6.49	113.39	108.20
1	AA	434	G	C5-C6-O6	-6.49	124.71	128.60
1	AA	1416	G	C5-C6-O6	-6.49	124.71	128.60
53	B5	92	G	C5-C6-O6	-6.49	124.71	128.60
53	B5	311	U	O4'-C1'-N1	6.49	113.39	108.20
53	B5	1061	A	O4'-C1'-N9	6.49	113.39	108.20
53	B5	1829	G	C5-C6-O6	-6.49	124.71	128.60
53	B5	1925	U	O4'-C1'-N1	6.49	113.39	108.20
1	AA	1080	G	C5-C6-O6	-6.48	124.71	128.60
53	B5	513	G	N1-C6-O6	6.48	123.79	119.90
53	B5	1174	G	O4'-C1'-N9	6.48	113.39	108.20
53	B5	1486	G	N1-C6-O6	6.48	123.79	119.90
53	B5	1727	G	N1-C6-O6	6.48	123.79	119.90
1	AA	79	C	O4'-C1'-N1	6.48	113.38	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
52	B4	7	U	O4'-C1'-N1	6.48	113.39	108.20
53	B5	1479	U	O4'-C1'-N1	6.48	113.38	108.20
53	B5	1906	G	O4'-C1'-N9	6.48	113.38	108.20
53	B5	2349	U	O4'-C1'-N1	6.48	113.38	108.20
1	AA	45	U	O4'-C1'-N1	6.48	113.38	108.20
1	AA	375	U	O4'-C1'-N1	6.48	113.38	108.20
1	AA	565	C	O4'-C1'-N1	6.48	113.38	108.20
1	AA	1393	U	O4'-C1'-N1	6.48	113.38	108.20
53	B5	1732	U	O4'-C1'-N1	6.48	113.38	108.20
53	B5	3194	C	O4'-C1'-N1	6.48	113.38	108.20
1	AA	1011	U	O4'-C1'-N1	6.48	113.38	108.20
51	B3	33	U	C2-N1-C1'	6.48	125.47	117.70
53	B5	98	G	N1-C6-O6	6.48	123.79	119.90
53	B5	440	A	O4'-C1'-N9	6.48	113.38	108.20
53	B5	1008	U	O4'-C1'-N1	6.48	113.38	108.20
53	B5	1060	U	O4'-C1'-N1	6.48	113.38	108.20
53	B5	2376	G	N1-C6-O6	6.48	123.79	119.90
53	B5	3222	U	O4'-C1'-N1	6.48	113.38	108.20
1	AA	474	A	O4'-C1'-N9	6.48	113.38	108.20
53	B5	1125	U	O4'-C1'-N1	6.48	113.38	108.20
53	B5	1168	U	O4'-C1'-N1	6.48	113.38	108.20
53	B5	2663	G	N1-C6-O6	6.48	123.79	119.90
53	B5	274	G	C5-C6-O6	-6.47	124.72	128.60
53	B5	889	U	O4'-C1'-N1	6.47	113.38	108.20
53	B5	941	G	N1-C6-O6	6.47	123.78	119.90
53	B5	1368	U	O4'-C1'-N1	6.47	113.38	108.20
1	AA	122	U	O4'-C1'-N1	6.47	113.38	108.20
51	B3	3	U	O4'-C1'-N1	6.47	113.38	108.20
51	B3	14	U	O4'-C1'-N1	6.47	113.38	108.20
53	B5	162	G	P-O3'-C3'	6.47	127.47	119.70
53	B5	1627	U	O4'-C1'-N1	6.47	113.38	108.20
53	B5	2168	A	O4'-C1'-N9	6.47	113.38	108.20
53	B5	2428	U	O4'-C1'-N1	6.47	113.38	108.20
53	B5	2663	G	O4'-C1'-N9	6.47	113.38	108.20
1	AA	688	G	C5-C6-O6	-6.47	124.72	128.60
1	AA	1150	G	O4'-C1'-N9	6.47	113.38	108.20
53	B5	75	G	O4'-C1'-N9	6.47	113.38	108.20
53	B5	495	G	N1-C6-O6	6.47	123.78	119.90
53	B5	1030	A	O4'-C1'-N9	6.47	113.38	108.20
53	B5	3254	G	N1-C6-O6	6.47	123.78	119.90
53	B5	3369	G	C5-C6-O6	-6.47	124.72	128.60
53	B5	72	C	O4'-C1'-N1	6.47	113.37	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	1617	G	N1-C6-O6	6.47	123.78	119.90
53	B5	2572	C	O4'-C1'-N1	6.47	113.37	108.20
53	B5	3016	A	O4'-C1'-N9	6.47	113.38	108.20
1	AA	199	G	N1-C6-O6	6.47	123.78	119.90
1	AA	246	G	O4'-C1'-N9	6.47	113.37	108.20
1	AA	277	U	O4'-C1'-N1	6.47	113.37	108.20
53	B5	865	U	O4'-C1'-N1	6.47	113.37	108.20
53	B5	2966	G	C5-C6-O6	-6.47	124.72	128.60
53	B5	3327	G	O4'-C1'-N9	6.47	113.37	108.20
53	B5	300	G	C5-C6-O6	-6.46	124.72	128.60
53	B5	558	U	O4'-C1'-N1	6.46	113.37	108.20
53	B5	2451	G	P-O3'-C3'	6.46	127.46	119.70
53	B5	2768	U	O4'-C1'-N1	6.46	113.37	108.20
1	AA	1322	A	O4'-C1'-N9	6.46	113.37	108.20
1	AA	1584	A	C8-N9-C4	-6.46	103.22	105.80
1	AA	1783	U	O4'-C1'-N1	6.46	113.37	108.20
52	B4	61	A	O4'-C1'-N9	6.46	113.37	108.20
52	B4	145	U	O4'-C1'-N1	6.46	113.37	108.20
53	B5	553	U	O4'-C1'-N1	6.46	113.37	108.20
53	B5	2684	C	O4'-C1'-N1	6.46	113.37	108.20
1	AA	656	G	N1-C6-O6	6.46	123.78	119.90
19	A7	15	G	N9-C1'-C2'	-6.46	104.89	112.00
19	A7	30	G	C8-N9-C4	-6.46	103.81	106.40
53	B5	2251	G	O4'-C1'-N9	6.46	113.37	108.20
1	AA	766	U	O4'-C1'-N1	6.46	113.37	108.20
53	B5	894	G	N1-C6-O6	6.46	123.78	119.90
1	AA	174	U	C2-N1-C1'	6.46	125.45	117.70
1	AA	232	U	O4'-C1'-N1	6.46	113.37	108.20
19	A7	27	C	C5'-C4'-O4'	6.46	116.85	109.10
19	A7	76	A	O4'-C1'-N9	6.46	113.37	108.20
53	B5	413	U	O4'-C1'-N1	6.46	113.37	108.20
1	AA	509	G	N1-C6-O6	6.46	123.77	119.90
1	AA	801	G	N1-C6-O6	6.46	123.77	119.90
1	AA	1229	G	N1-C6-O6	6.46	123.77	119.90
49	BY	47	TYR	CB-CG-CD2	6.46	124.87	121.00
1	AA	479	C	O4'-C1'-N1	6.45	113.36	108.20
53	B5	1674	G	N1-C6-O6	6.45	123.77	119.90
1	AA	231	U	O4'-C1'-N1	6.45	113.36	108.20
1	AA	839	U	O4'-C1'-N1	6.45	113.36	108.20
53	B5	541	U	O4'-C1'-N1	6.45	113.36	108.20
53	B5	2103	U	O4'-C1'-N1	6.45	113.36	108.20
1	AA	1724	G	N1-C6-O6	6.45	123.77	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
19	A7	43	G	N1-C6-O6	-6.45	116.03	119.90
53	B5	2530	G	O4'-C1'-N9	6.45	113.36	108.20
1	AA	613	G	C5-C6-O6	-6.45	124.73	128.60
1	AA	760	A	O4'-C1'-N9	6.45	113.36	108.20
1	AA	902	U	O4'-C1'-N1	6.45	113.36	108.20
53	B5	1733	G	C5-C6-O6	-6.45	124.73	128.60
1	AA	238	U	O4'-C1'-N1	6.45	113.36	108.20
53	B5	1560	G	N1-C6-O6	6.45	123.77	119.90
53	B5	3366	G	C5-C6-O6	-6.45	124.73	128.60
53	B5	1445	U	O4'-C1'-N1	6.44	113.36	108.20
53	B5	713	U	O4'-C1'-N1	6.44	113.35	108.20
53	B5	2477	G	C5-C6-O6	-6.44	124.73	128.60
53	B5	3066	U	O4'-C1'-N1	6.44	113.35	108.20
53	B5	1430	U	O4'-C1'-N1	6.44	113.35	108.20
53	B5	1442	U	O4'-C1'-N1	6.44	113.35	108.20
53	B5	2974	U	O4'-C1'-N1	6.44	113.35	108.20
53	B5	3371	G	C5-C6-O6	-6.44	124.73	128.60
1	AA	884	G	N1-C6-O6	6.44	123.76	119.90
1	AA	1511	G	C5-C6-O6	-6.44	124.74	128.60
1	AA	882	U	O4'-C1'-N1	6.44	113.35	108.20
1	AA	1652	G	N1-C6-O6	6.44	123.76	119.90
53	B5	129	U	O4'-C1'-N1	6.44	113.35	108.20
53	B5	1853	U	O4'-C1'-N1	6.44	113.35	108.20
1	AA	258	C	O4'-C1'-N1	6.44	113.35	108.20
1	AA	1674	U	O4'-C1'-N1	6.43	113.35	108.20
53	B5	411	U	O4'-C1'-N1	6.43	113.35	108.20
53	B5	1109	U	O4'-C1'-N1	6.43	113.35	108.20
53	B5	1751	G	O4'-C1'-N9	6.43	113.35	108.20
53	B5	3069	G	N1-C6-O6	6.43	123.76	119.90
1	AA	1353	U	O4'-C1'-N1	6.43	113.35	108.20
52	B4	143	U	O4'-C1'-N1	6.43	113.35	108.20
53	B5	721	G	N1-C6-O6	6.43	123.76	119.90
53	B5	769	G	C5-C6-O6	-6.43	124.74	128.60
53	B5	802	C	O4'-C1'-N1	6.43	113.34	108.20
53	B5	1029	G	N1-C6-O6	6.43	123.76	119.90
1	AA	1138	G	N1-C6-O6	6.43	123.76	119.90
53	B5	787	G	C5-C6-O6	-6.43	124.74	128.60
53	B5	1611	G	O4'-C1'-N9	6.43	113.34	108.20
1	AA	675	U	O4'-C1'-N1	6.43	113.34	108.20
1	AA	1151	G	O4'-C1'-N9	6.43	113.34	108.20
53	B5	128	G	O4'-C1'-N9	6.43	113.34	108.20
1	AA	1361	G	N1-C6-O6	6.43	123.75	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
52	B4	101	U	P-O3'-C3'	6.43	127.41	119.70
53	B5	123	A	O4'-C1'-N9	6.43	113.34	108.20
53	B5	1250	G	P-O3'-C3'	6.43	127.41	119.70
53	B5	1345	G	N1-C6-O6	6.43	123.76	119.90
53	B5	3081	C	O4'-C1'-N1	6.43	113.34	108.20
1	AA	465	G	C5-C6-O6	-6.42	124.75	128.60
1	AA	1560	G	N1-C6-O6	6.42	123.75	119.90
19	A7	33	U	N1-C2-N3	6.42	118.75	114.90
53	B5	1473	G	N1-C6-O6	6.42	123.75	119.90
53	B5	1717	U	O4'-C1'-N1	6.42	113.34	108.20
53	B5	2797	C	O4'-C1'-N1	6.42	113.34	108.20
1	AA	843	U	O4'-C1'-N1	6.42	113.34	108.20
1	AA	256	A	O4'-C1'-N9	6.42	113.34	108.20
1	AA	603	U	O4'-C1'-N1	6.42	113.34	108.20
1	AA	718	U	O4'-C1'-N1	6.42	113.34	108.20
1	AA	1739	U	O4'-C1'-N1	6.42	113.34	108.20
51	B3	13	A	O4'-C1'-N9	6.42	113.34	108.20
53	B5	243	G	O4'-C1'-N9	6.42	113.34	108.20
53	B5	988	U	O4'-C1'-N1	6.42	113.34	108.20
53	B5	2157	G	N1-C6-O6	6.42	123.75	119.90
53	B5	282	G	N1-C6-O6	6.42	123.75	119.90
1	AA	111	U	O4'-C1'-N1	6.42	113.34	108.20
1	AA	363	G	N1-C6-O6	6.42	123.75	119.90
1	AA	374	U	O4'-C1'-N1	6.42	113.33	108.20
53	B5	626	U	O4'-C1'-N1	6.42	113.33	108.20
53	B5	3189	G	N1-C6-O6	6.42	123.75	119.90
53	B5	3333	G	C5-C6-O6	-6.42	124.75	128.60
1	AA	58	U	O4'-C1'-N1	6.42	113.33	108.20
51	B3	81	U	O4'-C1'-N1	6.42	113.33	108.20
53	B5	688	G	N1-C6-O6	6.42	123.75	119.90
53	B5	772	U	O4'-C1'-N1	6.42	113.33	108.20
1	AA	155	U	O4'-C1'-N1	6.42	113.33	108.20
53	B5	2610	G	O4'-C1'-N9	6.42	113.33	108.20
1	AA	611	U	O4'-C1'-N1	6.41	113.33	108.20
1	AA	1386	C	C6-N1-C1'	-6.41	113.10	120.80
53	B5	444	U	O4'-C1'-N1	6.41	113.33	108.20
53	B5	652	G	N1-C6-O6	6.41	123.75	119.90
53	B5	852	U	O4'-C1'-N1	6.41	113.33	108.20
53	B5	1395	G	C5-C6-O6	-6.41	124.75	128.60
1	AA	1791	G	N1-C6-O6	6.41	123.75	119.90
53	B5	2509	U	O4'-C1'-N1	6.41	113.33	108.20
1	AA	318	U	O4'-C1'-N1	6.41	113.33	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	1616	U	O4'-C1'-N1	6.41	113.33	108.20
53	B5	2124	G	O4'-C1'-N9	6.41	113.33	108.20
1	AA	719	U	O4'-C1'-N1	6.41	113.33	108.20
53	B5	441	U	O4'-C1'-N1	6.41	113.33	108.20
53	B5	750	G	N1-C6-O6	6.41	123.75	119.90
53	B5	2787	G	N1-C6-O6	6.41	123.75	119.90
53	B5	2842	U	O4'-C1'-N1	6.41	113.33	108.20
1	AA	797	G	O4'-C4'-C3'	-6.41	97.59	104.00
53	B5	1592	G	C5-C6-O6	-6.41	124.76	128.60
53	B5	2994	A	O4'-C1'-N9	6.41	113.33	108.20
1	AA	1736	U	O4'-C1'-N1	6.41	113.32	108.20
53	B5	183	G	N1-C6-O6	6.41	123.74	119.90
53	B5	449	U	O4'-C1'-N1	6.41	113.32	108.20
53	B5	1077	U	O4'-C1'-N1	6.41	113.33	108.20
53	B5	2659	G	C5-C6-O6	-6.41	124.76	128.60
53	B5	2698	G	N1-C6-O6	6.41	123.74	119.90
53	B5	470	G	O4'-C1'-N9	6.40	113.32	108.20
53	B5	1095	U	O4'-C1'-N1	6.40	113.32	108.20
1	AA	452	A	O4'-C1'-N9	6.40	113.32	108.20
1	AA	1619	U	O4'-C1'-N1	6.40	113.32	108.20
53	B5	584	G	N1-C6-O6	6.40	123.74	119.90
53	B5	643	U	O4'-C1'-N1	6.40	113.32	108.20
53	B5	1246	G	C5-C6-O6	-6.40	124.76	128.60
53	B5	1577	G	N1-C6-O6	6.40	123.74	119.90
53	B5	2694	A	O4'-C1'-N9	6.40	113.32	108.20
53	B5	2814	G	N1-C6-O6	6.40	123.74	119.90
53	B5	3348	G	N1-C6-O6	6.40	123.74	119.90
53	B5	32	U	O4'-C1'-N1	6.40	113.32	108.20
53	B5	162	G	O4'-C1'-N9	6.40	113.32	108.20
53	B5	2793	G	C5-C6-O6	-6.40	124.76	128.60
1	AA	912	G	N1-C6-O6	6.40	123.74	119.90
53	B5	2918	G	O4'-C1'-N9	6.40	113.32	108.20
23	B8	18	TYR	CB-CG-CD1	6.40	124.84	121.00
1	AA	459	G	O4'-C1'-N9	6.40	113.32	108.20
1	AA	1446	G	C5-C6-O6	-6.40	124.76	128.60
53	B5	1116	G	C5-C6-O6	-6.40	124.76	128.60
53	B5	172	G	O4'-C1'-N9	6.39	113.31	108.20
53	B5	669	U	O4'-C1'-N1	6.39	113.31	108.20
53	B5	1295	G	O4'-C1'-N9	6.39	113.31	108.20
53	B5	1624	G	C5-C6-O6	-6.39	124.76	128.60
53	B5	2442	G	O4'-C1'-N9	6.39	113.31	108.20
1	AA	48	G	N1-C6-O6	6.39	123.74	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1389	U	O4'-C1'-N1	6.39	113.31	108.20
1	AA	81	G	N1-C6-O6	6.39	123.73	119.90
1	AA	642	G	N1-C6-O6	6.39	123.73	119.90
53	B5	548	G	N1-C6-O6	6.39	123.73	119.90
53	B5	2122	G	O4'-C1'-N9	6.39	113.31	108.20
53	B5	2598	G	O4'-C1'-N9	6.39	113.31	108.20
53	B5	3202	G	N1-C6-O6	6.39	123.73	119.90
53	B5	1576	G	O4'-C1'-N9	6.39	113.31	108.20
53	B5	1689	U	O4'-C1'-N1	6.39	113.31	108.20
53	B5	3260	G	O4'-C1'-N9	6.39	113.31	108.20
51	B3	112	G	O4'-C1'-N9	6.39	113.31	108.20
53	B5	326	U	O4'-C1'-N1	6.39	113.31	108.20
53	B5	1422	G	C5-C6-O6	-6.39	124.77	128.60
53	B5	3101	G	C5-C6-O6	-6.39	124.77	128.60
1	AA	163	G	N1-C6-O6	6.39	123.73	119.90
53	B5	158	G	C5-C6-O6	-6.39	124.77	128.60
53	B5	2199	G	O4'-C1'-N9	6.39	113.31	108.20
53	B5	2236	G	N1-C6-O6	6.39	123.73	119.90
53	B5	2310	U	O4'-C1'-N1	6.39	113.31	108.20
53	B5	2908	G	N1-C6-O6	6.39	123.73	119.90
1	AA	487	G	N1-C6-O6	6.38	123.73	119.90
1	AA	1786	G	C5-C6-O6	-6.38	124.77	128.60
53	B5	297	G	C5-C6-O6	-6.38	124.77	128.60
53	B5	1752	A	O4'-C1'-N9	6.38	113.31	108.20
52	B4	5	U	O4'-C1'-N1	6.38	113.31	108.20
53	B5	1385	C	O4'-C1'-N1	6.38	113.31	108.20
1	AA	1054	U	O4'-C1'-N1	6.38	113.31	108.20
1	AA	1124	G	N1-C6-O6	6.38	123.73	119.90
51	B3	56	G	N1-C6-O6	6.38	123.73	119.90
53	B5	2923	U	O4'-C1'-N1	6.38	113.31	108.20
53	B5	3022	G	C5-C6-O6	-6.38	124.77	128.60
1	AA	27	U	O4'-C1'-N1	6.38	113.30	108.20
1	AA	63	G	N1-C6-O6	6.38	123.73	119.90
1	AA	1023	U	O4'-C1'-N1	6.38	113.30	108.20
53	B5	2423	U	O4'-C1'-N1	6.38	113.30	108.20
53	B5	2790	A	O4'-C1'-N9	6.38	113.30	108.20
1	AA	540	G	O4'-C1'-N9	6.38	113.30	108.20
53	B5	910	G	C5-C6-O6	-6.38	124.77	128.60
53	B5	739	G	N1-C6-O6	6.38	123.73	119.90
53	B5	879	U	O4'-C1'-N1	6.38	113.30	108.20
53	B5	2891	U	O4'-C1'-N1	6.38	113.30	108.20
53	B5	2964	G	C5-C6-O6	-6.38	124.78	128.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	3111	U	O4'-C1'-N1	6.38	113.30	108.20
1	AA	652	G	N1-C6-O6	6.38	123.72	119.90
1	AA	1247	U	O4'-C1'-N1	6.38	113.30	108.20
53	B5	3077	A	O4'-C1'-N9	6.38	113.30	108.20
1	AA	200	A	O4'-C1'-N9	6.37	113.30	108.20
52	B4	103	G	C5-C6-O6	-6.37	124.78	128.60
53	B5	468	G	O4'-C1'-N9	6.37	113.30	108.20
53	B5	1928	G	C5-C6-O6	-6.37	124.78	128.60
53	B5	2725	U	O4'-C1'-N1	6.37	113.30	108.20
19	A7	22	G	P-O3'-C3'	6.37	127.35	119.70
19	A7	42	G	C5-N7-C8	-6.37	101.11	104.30
53	B5	400	G	N1-C6-O6	6.37	123.72	119.90
53	B5	2683	U	O4'-C1'-N1	6.37	113.30	108.20
1	AA	1241	G	C5-C6-O6	-6.37	124.78	128.60
1	AA	1575	A	O4'-C1'-N9	6.37	113.30	108.20
53	B5	317	A	P-O3'-C3'	6.37	127.34	119.70
53	B5	1289	G	N1-C6-O6	6.37	123.72	119.90
53	B5	1871	U	O4'-C1'-N1	6.37	113.30	108.20
1	AA	1533	U	O4'-C1'-N1	6.37	113.29	108.20
51	B3	2	G	N1-C6-O6	6.37	123.72	119.90
53	B5	21	G	N1-C6-O6	6.37	123.72	119.90
53	B5	1080	A	O4'-C1'-N9	6.37	113.29	108.20
53	B5	1467	A	O4'-C1'-N9	6.37	113.29	108.20
53	B5	3386	G	C5-C6-O6	-6.36	124.78	128.60
1	AA	1267	G	N1-C6-O6	6.36	123.72	119.90
19	A7	68	U	C4'-C3'-C2'	-6.36	96.24	102.60
53	B5	1174	G	N1-C6-O6	6.36	123.72	119.90
53	B5	3158	G	N1-C6-O6	6.36	123.72	119.90
1	AA	159	U	O4'-C1'-N1	6.36	113.29	108.20
53	B5	1720	U	O4'-C1'-N1	6.36	113.29	108.20
1	AA	676	G	N1-C6-O6	6.36	123.71	119.90
1	AA	1142	U	O4'-C1'-N1	6.36	113.29	108.20
53	B5	3356	G	N1-C6-O6	6.36	123.71	119.90
53	B5	1747	G	O4'-C1'-N9	6.36	113.28	108.20
53	B5	2882	U	O4'-C1'-N1	6.36	113.28	108.20
53	B5	3196	U	O4'-C1'-N1	6.36	113.28	108.20
53	B5	3212	U	O4'-C1'-N1	6.36	113.28	108.20
1	AA	1711	G	N1-C6-O6	6.35	123.71	119.90
52	B4	26	U	O4'-C1'-N1	6.35	113.28	108.20
53	B5	56	G	O4'-C1'-N9	6.35	113.28	108.20
53	B5	1021	G	N1-C6-O6	6.35	123.71	119.90
53	B5	1189	C	O4'-C1'-N1	6.35	113.28	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	2833	A	C5-C6-N6	-6.35	118.62	123.70
53	B5	3241	G	C5-C6-O6	-6.35	124.79	128.60
1	AA	142	G	N1-C6-O6	6.35	123.71	119.90
1	AA	1539	G	C5-C6-O6	-6.35	124.79	128.60
19	A7	68	U	N1-C2-N3	6.35	118.71	114.90
53	B5	640	U	O4'-C1'-N1	6.35	113.28	108.20
53	B5	658	G	N1-C6-O6	6.35	123.71	119.90
53	B5	3384	U	O4'-C1'-N1	6.35	113.28	108.20
1	AA	1208	G	O4'-C1'-N9	6.35	113.28	108.20
1	AA	1415	G	N1-C6-O6	6.35	123.71	119.90
53	B5	2481	G	C5-C6-O6	-6.35	124.79	128.60
53	B5	2980	U	O4'-C1'-N1	6.35	113.28	108.20
53	B5	3246	G	N1-C6-O6	6.35	123.71	119.90
1	AA	1476	G	N1-C6-O6	6.34	123.71	119.90
19	A7	59	U	N3-C4-C5	-6.34	110.79	114.60
53	B5	1666	G	C5-C6-O6	-6.34	124.79	128.60
53	B5	2385	G	C5-C6-O6	-6.34	124.79	128.60
53	B5	3290	G	C5-C6-O6	-6.34	124.79	128.60
1	AA	1250	U	O4'-C1'-N1	6.34	113.28	108.20
1	AA	1375	U	O4'-C1'-N1	6.34	113.27	108.20
53	B5	2332	A	O4'-C1'-N9	6.34	113.27	108.20
53	B5	2355	G	C5-C6-O6	-6.34	124.79	128.60
1	AA	372	G	O4'-C1'-N9	6.34	113.27	108.20
1	AA	1166	G	N1-C6-O6	6.34	123.70	119.90
1	AA	1348	G	O4'-C1'-N9	6.34	113.27	108.20
53	B5	2342	U	O4'-C1'-N1	6.34	113.27	108.20
1	AA	291	G	N1-C6-O6	6.34	123.70	119.90
53	B5	2417	U	O4'-C1'-N1	6.34	113.27	108.20
53	B5	2498	U	O4'-C1'-N1	6.34	113.27	108.20
1	AA	595	G	O4'-C1'-N9	6.34	113.27	108.20
51	B3	21	G	C5-C6-O6	-6.34	124.80	128.60
52	B4	58	G	P-O3'-C3'	6.34	127.31	119.70
19	A7	66	A	C4-C5-C6	-6.34	113.83	117.00
53	B5	975	G	N1-C6-O6	6.34	123.70	119.90
53	B5	1677	G	N1-C6-O6	6.34	123.70	119.90
53	B5	3352	U	O4'-C1'-N1	6.34	113.27	108.20
53	B5	333	G	O4'-C1'-N9	6.33	113.27	108.20
1	AA	379	U	O4'-C1'-N1	6.33	113.27	108.20
1	AA	806	A	O4'-C1'-N9	6.33	113.27	108.20
1	AA	908	U	O4'-C1'-N1	6.33	113.27	108.20
1	AA	916	U	O4'-C1'-N1	6.33	113.27	108.20
53	B5	461	U	O4'-C1'-N1	6.33	113.27	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	760	G	N1-C6-O6	6.33	123.70	119.90
53	B5	1115	G	N1-C6-O6	6.33	123.70	119.90
53	B5	3325	G	O4'-C1'-N9	6.33	113.27	108.20
1	AA	648	G	C5-C6-O6	-6.33	124.80	128.60
19	A7	57	G	N1-C6-O6	-6.33	116.10	119.90
52	B4	10	A	O4'-C1'-N9	6.33	113.27	108.20
52	B4	42	G	N1-C6-O6	6.33	123.70	119.90
53	B5	281	G	C5-C6-O6	-6.33	124.80	128.60
53	B5	390	G	N1-C6-O6	6.33	123.70	119.90
53	B5	495	G	O4'-C1'-N9	6.33	113.27	108.20
53	B5	2574	G	O4'-C1'-N9	6.33	113.27	108.20
1	AA	1016	U	O4'-C1'-N1	6.33	113.26	108.20
19	A7	19	G	N3-C2-N2	-6.33	115.47	119.90
53	B5	164	A	O4'-C1'-N9	6.33	113.26	108.20
53	B5	2247	G	O4'-C1'-N9	6.33	113.26	108.20
53	B5	829	U	O4'-C1'-N1	6.33	113.26	108.20
53	B5	2864	A	P-O3'-C3'	6.33	127.29	119.70
53	B5	1498	A	O4'-C1'-N9	6.33	113.26	108.20
53	B5	564	G	O4'-C1'-N9	6.33	113.26	108.20
53	B5	878	G	O4'-C1'-N9	6.33	113.26	108.20
53	B5	1412	G	C5-C6-O6	-6.33	124.80	128.60
53	B5	1780	G	N1-C6-O6	6.33	123.70	119.90
1	AA	695	U	O4'-C1'-N1	6.32	113.26	108.20
1	AA	1001	G	C5-C6-O6	-6.32	124.81	128.60
1	AA	1583	U	O4'-C1'-N1	6.32	113.26	108.20
51	B3	82	A	O4'-C1'-N9	6.32	113.26	108.20
53	B5	87	U	O4'-C1'-N1	6.32	113.26	108.20
53	B5	314	U	O4'-C1'-N1	6.32	113.26	108.20
1	AA	1109	G	N1-C6-O6	6.32	123.69	119.90
1	AA	1433	A	O4'-C1'-N9	6.32	113.26	108.20
52	B4	98	U	O4'-C1'-N1	6.32	113.26	108.20
53	B5	924	G	N1-C6-O6	6.32	123.69	119.90
53	B5	2336	U	O4'-C1'-N1	6.32	113.26	108.20
1	AA	1259	G	N1-C6-O6	6.32	123.69	119.90
1	AA	1609	A	O4'-C1'-N9	6.32	113.26	108.20
1	AA	1267	G	O4'-C1'-N9	6.32	113.25	108.20
53	B5	1024	G	P-O3'-C3'	6.32	127.28	119.70
53	B5	2943	G	C5-C6-O6	-6.32	124.81	128.60
1	AA	368	U	O4'-C1'-N1	6.32	113.25	108.20
1	AA	858	G	C2-N3-C4	6.32	115.06	111.90
1	AA	1656	G	N1-C6-O6	6.32	123.69	119.90
53	B5	1718	G	P-O3'-C3'	6.32	127.28	119.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	2317	A	O4'-C1'-N9	6.32	113.25	108.20
53	B5	3156	U	O4'-C1'-N1	6.32	113.25	108.20
53	B5	3390	G	C5-C6-O6	-6.32	124.81	128.60
53	B5	1393	A	O4'-C1'-N9	6.31	113.25	108.20
1	AA	564	G	C5-C6-O6	-6.31	124.81	128.60
52	B4	144	G	C5-C6-O6	-6.31	124.81	128.60
53	B5	512	U	O4'-C1'-N1	6.31	113.25	108.20
53	B5	2250	G	O4'-C1'-N9	6.31	113.25	108.20
53	B5	1927	G	O4'-C1'-N9	6.31	113.25	108.20
1	AA	960	U	O4'-C1'-N1	6.31	113.25	108.20
1	AA	1427	U	O4'-C1'-N1	6.31	113.25	108.20
1	AA	1759	U	O4'-C1'-N1	6.31	113.25	108.20
53	B5	1056	U	O4'-C1'-N1	6.31	113.25	108.20
1	AA	548	G	O4'-C1'-N9	6.31	113.25	108.20
1	AA	1550	U	C4'-C3'-C2'	6.31	108.91	102.60
52	B4	39	G	C5-C6-O6	-6.31	124.81	128.60
53	B5	2	U	O4'-C1'-N1	6.31	113.25	108.20
53	B5	853	G	O4'-C1'-N9	6.31	113.25	108.20
53	B5	1082	U	O4'-C1'-N1	6.31	113.25	108.20
53	B5	1134	G	O4'-C1'-N9	6.31	113.25	108.20
53	B5	1233	G	N1-C6-O6	6.31	123.69	119.90
53	B5	2534	G	O4'-C1'-N9	6.31	113.25	108.20
53	B5	2631	U	O4'-C1'-N1	6.31	113.25	108.20
53	B5	2787	G	O4'-C1'-N9	6.31	113.25	108.20
1	AA	52	U	O4'-C1'-N1	6.30	113.24	108.20
1	AA	643	G	N1-C6-O6	6.30	123.68	119.90
38	BN	4	TYR	CB-CG-CD1	6.30	124.78	121.00
53	B5	731	U	O4'-C1'-N1	6.30	113.24	108.20
53	B5	1036	A	O4'-C1'-N9	6.30	113.24	108.20
53	B5	1492	G	C5-C6-O6	-6.30	124.82	128.60
53	B5	1905	G	C5-C6-O6	-6.30	124.82	128.60
53	B5	1552	G	N1-C6-O6	6.30	123.68	119.90
1	AA	482	U	O4'-C1'-N1	6.30	113.24	108.20
1	AA	1732	U	O4'-C1'-N1	6.30	113.24	108.20
53	B5	244	G	N1-C6-O6	6.30	123.68	119.90
53	B5	3199	G	P-O3'-C3'	6.30	127.26	119.70
53	B5	3242	G	C5-C6-O6	-6.30	124.82	128.60
53	B5	3267	U	O4'-C1'-N1	6.30	113.24	108.20
1	AA	1520	U	O5'-P-OP1	-6.30	100.03	105.70
53	B5	20	A	O4'-C1'-N9	6.30	113.24	108.20
53	B5	355	A	O4'-C1'-N9	6.30	113.24	108.20
53	B5	2584	G	N1-C6-O6	6.30	123.68	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	2907	G	O4'-C1'-N9	6.30	113.24	108.20
53	B5	673	U	O4'-C1'-N1	6.30	113.24	108.20
53	B5	1476	G	C5-C6-O6	-6.30	124.82	128.60
53	B5	2325	G	C5-C6-O6	-6.30	124.82	128.60
53	B5	2403	G	N1-C6-O6	6.30	123.68	119.90
53	B5	2968	G	N1-C6-O6	6.30	123.68	119.90
1	AA	327	U	O4'-C1'-N1	6.29	113.24	108.20
1	AA	1695	G	O4'-C1'-N9	6.29	113.24	108.20
53	B5	295	A	O4'-C1'-N9	6.29	113.24	108.20
1	AA	167	U	O4'-C1'-N1	6.29	113.23	108.20
1	AA	1321	G	N1-C6-O6	6.29	123.68	119.90
1	AA	1523	A	OP1-P-OP2	-6.29	110.16	119.60
51	B3	43	U	P-O3'-C3'	6.29	127.25	119.70
51	B3	109	G	O4'-C1'-N9	6.29	113.23	108.20
52	B4	82	U	O4'-C1'-N1	6.29	113.23	108.20
52	B4	128	U	O4'-C1'-N1	6.29	113.23	108.20
53	B5	104	G	O4'-C1'-N9	6.29	113.23	108.20
53	B5	705	A	C5-C6-N6	-6.29	118.67	123.70
53	B5	1331	U	O4'-C1'-N1	6.29	113.23	108.20
1	AA	669	G	N1-C6-O6	6.29	123.67	119.90
53	B5	55	G	N1-C6-O6	6.29	123.67	119.90
53	B5	413	U	P-O3'-C3'	6.29	127.25	119.70
53	B5	1329	U	O4'-C1'-N1	6.29	113.23	108.20
53	B5	1475	A	O4'-C1'-N9	6.29	113.23	108.20
53	B5	1630	U	O4'-C1'-N1	6.29	113.23	108.20
53	B5	2769	A	O4'-C1'-N9	6.29	113.23	108.20
1	AA	77	U	O4'-C1'-N1	6.29	113.23	108.20
19	A7	60	C	C3'-C2'-C1'	-6.29	96.47	101.50
53	B5	1222	G	C5-C6-O6	-6.29	124.83	128.60
1	AA	738	G	N1-C6-O6	6.29	123.67	119.90
1	AA	1255	U	O4'-C1'-N1	6.29	113.23	108.20
1	AA	1462	G	O4'-C1'-N9	6.29	113.23	108.20
19	A7	2	C	N3-C4-C5	6.29	124.42	121.90
53	B5	426	G	N1-C6-O6	6.29	123.67	119.90
53	B5	1131	G	C5-C6-O6	-6.29	124.83	128.60
53	B5	1229	G	N1-C6-O6	6.29	123.67	119.90
1	AA	1332	U	O4'-C1'-N1	6.29	113.23	108.20
19	A7	71	G	C4-C5-N7	-6.29	108.28	110.80
53	B5	1493	G	C5-C6-O6	-6.29	124.83	128.60
1	AA	89	G	C5-C6-O6	-6.29	124.83	128.60
1	AA	980	U	O4'-C1'-N1	6.29	113.23	108.20
53	B5	754	G	O4'-C1'-N9	6.29	113.23	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	1114	U	O4'-C1'-N1	6.29	113.23	108.20
53	B5	1486	G	O4'-C1'-N9	6.29	113.23	108.20
53	B5	1886	A	O4'-C1'-N9	6.29	113.23	108.20
53	B5	3389	U	O4'-C1'-N1	6.29	113.23	108.20
1	AA	306	U	O4'-C1'-N1	6.28	113.23	108.20
1	AA	423	G	N1-C6-O6	6.28	123.67	119.90
1	AA	511	A	O4'-C1'-N9	6.28	113.23	108.20
1	AA	774	A	O4'-C1'-N9	6.28	113.23	108.20
1	AA	1189	A	O4'-C1'-N9	6.28	113.23	108.20
1	AA	1269	G	O4'-C1'-N9	6.28	113.23	108.20
53	B5	222	A	O4'-C1'-N9	6.28	113.23	108.20
53	B5	726	G	N1-C6-O6	6.28	123.67	119.90
1	AA	1112	U	O4'-C1'-N1	6.28	113.22	108.20
1	AA	1777	U	O4'-C1'-N1	6.28	113.23	108.20
53	B5	1842	A	O4'-C1'-N9	6.28	113.22	108.20
1	AA	390	G	O4'-C1'-N9	6.28	113.22	108.20
52	B4	64	U	O4'-C1'-N1	6.28	113.22	108.20
53	B5	25	U	O4'-C1'-N1	6.28	113.22	108.20
53	B5	1249	G	O4'-C1'-N9	6.28	113.22	108.20
53	B5	1577	G	O4'-C1'-N9	6.28	113.22	108.20
53	B5	3306	U	O4'-C1'-N1	6.28	113.22	108.20
1	AA	714	G	O4'-C1'-N9	6.28	113.22	108.20
1	AA	900	G	N1-C6-O6	6.28	123.67	119.90
51	B3	8	G	C5-C6-O6	-6.28	124.83	128.60
53	B5	3083	G	N1-C6-O6	6.28	123.67	119.90
1	AA	115	G	O4'-C1'-N9	6.28	113.22	108.20
1	AA	663	U	O4'-C1'-N1	6.28	113.22	108.20
1	AA	1213	A	O4'-C1'-N9	6.28	113.22	108.20
36	BL	141	ALA	N-CA-CB	6.28	118.89	110.10
53	B5	161	G	O4'-C1'-N9	6.28	113.22	108.20
53	B5	383	G	C5-C6-O6	-6.28	124.83	128.60
53	B5	716	G	C5-C6-O6	-6.28	124.83	128.60
53	B5	1134	G	N1-C6-O6	6.28	123.67	119.90
53	B5	1483	G	N1-C6-O6	6.28	123.67	119.90
53	B5	2871	G	N1-C6-O6	6.28	123.67	119.90
1	AA	20	G	O4'-C1'-N9	6.28	113.22	108.20
1	AA	585	A	O4'-C1'-N9	6.28	113.22	108.20
1	AA	1114	U	O4'-C1'-N1	6.28	113.22	108.20
1	AA	1176	G	N1-C6-O6	6.28	123.67	119.90
1	AA	1252	A	O4'-C1'-N9	6.28	113.22	108.20
19	A7	18	G	C4'-C3'-C2'	-6.28	96.33	102.60
53	B5	3147	G	N1-C6-O6	6.28	123.67	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
17	AR	285	ALA	C-N-CA	6.27	137.38	121.70
52	B4	13	A	O4'-C1'-N9	6.27	113.22	108.20
1	AA	983	G	O4'-C1'-N9	6.27	113.22	108.20
53	B5	1890	U	O4'-C1'-N1	6.27	113.22	108.20
1	AA	1165	U	O4'-C1'-N1	6.27	113.22	108.20
53	B5	371	G	N1-C6-O6	6.27	123.66	119.90
53	B5	603	A	O4'-C1'-N9	6.27	113.22	108.20
53	B5	722	G	N1-C6-O6	6.27	123.66	119.90
53	B5	885	U	O4'-C1'-N1	6.27	113.22	108.20
53	B5	2454	G	O4'-C1'-N9	6.27	113.22	108.20
53	B5	2472	U	O4'-C1'-N1	6.27	113.21	108.20
53	B5	2800	G	N1-C6-O6	6.27	123.66	119.90
53	B5	1377	G	O4'-C1'-N9	6.27	113.21	108.20
53	B5	2253	G	O4'-C1'-N9	6.27	113.21	108.20
53	B5	3274	G	O4'-C1'-N9	6.27	113.21	108.20
1	AA	201	G	N1-C6-O6	6.26	123.66	119.90
1	AA	252	U	O4'-C1'-N1	6.26	113.21	108.20
1	AA	714	G	N1-C6-O6	6.26	123.66	119.90
53	B5	2105	G	O4'-C1'-N9	6.26	113.21	108.20
53	B5	2672	G	N1-C6-O6	6.26	123.66	119.90
53	B5	1649	U	O4'-C1'-N1	6.26	113.21	108.20
53	B5	1770	G	C5-C6-O6	-6.26	124.84	128.60
53	B5	3146	G	N1-C6-O6	6.26	123.66	119.90
1	AA	392	G	O4'-C1'-N9	6.26	113.21	108.20
1	AA	802	G	O4'-C1'-N9	6.26	113.21	108.20
1	AA	896	U	O4'-C1'-N1	6.26	113.21	108.20
1	AA	919	U	O4'-C1'-N1	6.26	113.21	108.20
1	AA	1107	G	N1-C6-O6	6.26	123.66	119.90
1	AA	1307	U	O4'-C1'-N1	6.26	113.21	108.20
53	B5	3058	U	O4'-C1'-N1	6.26	113.21	108.20
1	AA	1048	G	N1-C6-O6	6.26	123.66	119.90
1	AA	1443	A	O4'-C1'-N9	6.26	113.21	108.20
1	AA	1693	G	O4'-C1'-N9	6.26	113.21	108.20
1	AA	823	G	O4'-C1'-N9	6.26	113.21	108.20
1	AA	1532	G	N1-C6-O6	6.26	123.66	119.90
53	B5	3179	U	O4'-C1'-N1	6.26	113.21	108.20
1	AA	53	G	C5-C6-O6	-6.26	124.85	128.60
42	BR	94	TYR	CB-CG-CD2	6.26	124.75	121.00
52	B4	116	G	O4'-C1'-N9	6.26	113.20	108.20
53	B5	685	G	O4'-C1'-N9	6.26	113.20	108.20
1	AA	143	G	N1-C6-O6	6.25	123.65	119.90
1	AA	609	U	O4'-C1'-N1	6.25	113.20	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1103	U	O4'-C1'-N1	6.25	113.20	108.20
1	AA	1628	U	O4'-C1'-N1	6.25	113.20	108.20
1	AA	239	C	O4'-C1'-N1	6.25	113.20	108.20
1	AA	1091	G	N1-C6-O6	6.25	123.65	119.90
1	AA	1288	G	O4'-C1'-N9	6.25	113.20	108.20
1	AA	1488	C	O4'-C1'-N1	6.25	113.20	108.20
53	B5	972	G	N1-C6-O6	6.25	123.65	119.90
53	B5	1892	G	N1-C6-O6	6.25	123.65	119.90
53	B5	2249	G	O4'-C1'-N9	6.25	113.20	108.20
1	AA	1147	G	N1-C6-O6	6.25	123.65	119.90
53	B5	3309	G	C5-C6-O6	-6.25	124.85	128.60
1	AA	33	U	O4'-C1'-N1	6.25	113.20	108.20
1	AA	348	U	O4'-C1'-N1	6.25	113.20	108.20
53	B5	2975	U	O4'-C1'-N1	6.25	113.20	108.20
53	B5	3055	U	O4'-C1'-N1	6.25	113.20	108.20
1	AA	549	G	N1-C6-O6	6.25	123.65	119.90
1	AA	857	U	C5'-C4'-O4'	6.25	116.59	109.10
18	AT	18	TYR	CB-CG-CD1	6.25	124.75	121.00
53	B5	302	U	O4'-C1'-N1	6.25	113.20	108.20
53	B5	2191	U	O4'-C1'-N1	6.25	113.20	108.20
1	AA	641	G	C5-C6-O6	-6.25	124.85	128.60
53	B5	748	U	O4'-C1'-N1	6.25	113.20	108.20
53	B5	1753	G	O4'-C1'-N9	6.25	113.20	108.20
1	AA	1100	U	O4'-C1'-N1	6.24	113.19	108.20
1	AA	1167	G	N1-C6-O6	6.24	123.65	119.90
1	AA	1183	U	O4'-C1'-N1	6.24	113.19	108.20
1	AA	1380	G	O4'-C1'-N9	6.24	113.19	108.20
52	B4	85	G	O4'-C1'-N9	6.24	113.19	108.20
53	B5	2579	G	C5-C6-O6	-6.24	124.85	128.60
1	AA	1299	U	O4'-C1'-N1	6.24	113.19	108.20
1	AA	685	A	C5-C6-N6	-6.24	118.71	123.70
53	B5	215	G	N1-C6-O6	6.24	123.64	119.90
53	B5	671	U	O4'-C1'-N1	6.24	113.19	108.20
53	B5	2218	G	C5-C6-O6	-6.24	124.86	128.60
53	B5	2863	G	N1-C6-O6	6.24	123.64	119.90
1	AA	458	G	N1-C6-O6	6.24	123.64	119.90
19	A7	22	G	C5-N7-C8	-6.24	101.18	104.30
52	B4	3	A	O4'-C1'-N9	6.24	113.19	108.20
53	B5	542	G	N1-C6-O6	6.24	123.64	119.90
53	B5	898	U	O4'-C1'-N1	6.24	113.19	108.20
53	B5	1670	C	O4'-C1'-N1	6.24	113.19	108.20
53	B5	1677	G	O4'-C1'-N9	6.24	113.19	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	2177	G	O4'-C1'-N9	6.24	113.19	108.20
53	B5	2221	G	C5-C6-O6	-6.24	124.86	128.60
53	B5	3283	U	O4'-C1'-N1	6.24	113.19	108.20
53	B5	171	G	O4'-C1'-N9	6.24	113.19	108.20
53	B5	2116	G	N1-C6-O6	6.24	123.64	119.90
1	AA	632	U	O4'-C1'-N1	6.24	113.19	108.20
1	AA	782	U	O4'-C1'-N1	6.24	113.19	108.20
1	AA	810	G	O4'-C1'-N9	6.24	113.19	108.20
1	AA	910	U	O4'-C1'-N1	6.24	113.19	108.20
53	B5	294	U	O4'-C1'-N1	6.24	113.19	108.20
53	B5	1634	G	O4'-C1'-N9	6.24	113.19	108.20
53	B5	2741	C	O4'-C1'-N1	6.24	113.19	108.20
53	B5	613	G	C5-C6-O6	-6.23	124.86	128.60
1	AA	1630	C	O5'-P-OP2	-6.23	100.09	105.70
53	B5	207	U	O4'-C1'-N1	6.23	113.19	108.20
53	B5	880	G	C5-C6-O6	-6.23	124.86	128.60
53	B5	1807	G	N1-C6-O6	6.23	123.64	119.90
53	B5	1828	A	O4'-C1'-N9	6.23	113.19	108.20
1	AA	1522	A	O4'-C1'-N9	-6.23	103.22	108.20
1	AA	1793	U	O4'-C1'-N1	6.23	113.19	108.20
19	A7	1	G	N3-C2-N2	-6.23	115.54	119.90
51	B3	109	G	N1-C6-O6	6.23	123.64	119.90
53	B5	536	U	O4'-C1'-N1	6.23	113.19	108.20
53	B5	1705	U	O4'-C1'-N1	6.23	113.19	108.20
53	B5	148	G	C5-C6-O6	-6.23	124.86	128.60
1	AA	1536	U	O4'-C1'-N1	6.23	113.18	108.20
1	AA	1734	G	O4'-C1'-N9	6.23	113.18	108.20
52	B4	58	G	O4'-C1'-N9	6.23	113.18	108.20
53	B5	231	G	N1-C6-O6	6.23	123.64	119.90
53	B5	341	G	N1-C6-O6	6.23	123.64	119.90
53	B5	2547	A	O4'-C1'-N9	6.23	113.18	108.20
53	B5	2662	G	O4'-C1'-N9	6.23	113.18	108.20
53	B5	3104	U	O4'-C1'-N1	6.23	113.18	108.20
1	AA	772	G	O4'-C1'-N9	6.23	113.18	108.20
53	B5	1747	G	C5-C6-O6	-6.23	124.86	128.60
53	B5	3076	C	O4'-C1'-N1	6.23	113.18	108.20
1	AA	1355	G	C5-C6-O6	-6.22	124.87	128.60
53	B5	979	G	O4'-C1'-N9	6.22	113.18	108.20
1	AA	952	G	C5-C6-O6	-6.22	124.87	128.60
1	AA	1257	G	N1-C6-O6	6.22	123.63	119.90
53	B5	992	G	C5-C6-O6	-6.22	124.87	128.60
53	B5	3149	G	O4'-C1'-N9	6.22	113.18	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	3213	G	C5-C6-O6	-6.22	124.87	128.60
53	B5	3318	G	C5-C6-O6	-6.22	124.87	128.60
1	AA	913	G	N1-C6-O6	6.22	123.63	119.90
53	B5	488	U	O4'-C1'-N1	6.22	113.18	108.20
53	B5	2759	U	O4'-C1'-N1	6.22	113.18	108.20
1	AA	418	G	O4'-C1'-N9	6.22	113.18	108.20
1	AA	1573	G	N1-C6-O6	6.22	123.63	119.90
52	B4	148	G	N1-C6-O6	6.22	123.63	119.90
53	B5	3332	U	O4'-C1'-N1	6.22	113.18	108.20
53	B5	518	G	N1-C6-O6	6.22	123.63	119.90
53	B5	2394	G	N1-C6-O6	6.22	123.63	119.90
17	AR	286	GLU	CB-CA-C	-6.21	97.97	110.40
51	B3	80	G	N1-C6-O6	6.21	123.63	119.90
53	B5	301	G	N1-C6-O6	6.21	123.63	119.90
53	B5	1450	G	C5-C6-O6	-6.21	124.87	128.60
53	B5	1830	G	C5-C6-O6	-6.21	124.87	128.60
53	B5	2393	G	C5-C6-O6	-6.21	124.87	128.60
53	B5	3230	G	O4'-C1'-N9	6.21	113.17	108.20
53	B5	2483	G	C5-C6-O6	-6.21	124.87	128.60
53	B5	3224	G	O4'-C1'-N9	6.21	113.17	108.20
29	BE	224	LYS	N-CA-CB	6.21	121.78	110.60
53	B5	414	U	O4'-C1'-N1	6.21	113.17	108.20
53	B5	2726	C	O4'-C1'-N1	6.21	113.17	108.20
53	B5	3297	U	O4'-C1'-N1	6.21	113.17	108.20
1	AA	364	G	N1-C6-O6	6.21	123.63	119.90
1	AA	1360	U	O4'-C1'-N1	6.21	113.17	108.20
19	A7	38	A	C6-C5-N7	6.21	136.65	132.30
53	B5	3035	A	O4'-C1'-N9	6.21	113.17	108.20
1	AA	1790	G	C5-C6-O6	-6.21	124.88	128.60
19	A7	33	U	C6-N1-C2	-6.21	117.28	121.00
53	B5	1087	G	P-O3'-C3'	6.21	127.15	119.70
53	B5	1777	U	C2-N1-C1'	6.21	125.15	117.70
1	AA	1117	U	O4'-C1'-N1	6.21	113.17	108.20
1	AA	1282	U	O4'-C1'-N1	6.21	113.17	108.20
53	B5	246	U	O4'-C1'-N1	6.21	113.17	108.20
53	B5	2319	U	O4'-C1'-N1	6.21	113.17	108.20
1	AA	1740	U	O4'-C1'-N1	6.20	113.16	108.20
19	A7	42	G	C6-C5-N7	-6.20	126.68	130.40
53	B5	624	G	N1-C6-O6	6.20	123.62	119.90
53	B5	1576	G	N1-C6-O6	6.20	123.62	119.90
53	B5	2624	G	C5-C6-O6	-6.20	124.88	128.60
1	AA	1349	G	N1-C6-O6	6.20	123.62	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	644	G	C5-C6-O6	-6.20	124.88	128.60
1	AA	1047	G	O4'-C1'-N9	6.20	113.16	108.20
19	A7	2	C	C5'-C4'-O4'	6.20	116.54	109.10
53	B5	511	G	N1-C6-O6	6.20	123.62	119.90
53	B5	901	G	O4'-C1'-N9	6.20	113.16	108.20
53	B5	1536	G	O4'-C1'-N9	6.20	113.16	108.20
53	B5	3220	G	O4'-C1'-N9	6.20	113.16	108.20
1	AA	774	A	P-O3'-C3'	-6.20	112.26	119.70
1	AA	1702	U	O4'-C1'-N1	6.20	113.16	108.20
53	B5	212	G	C5-C6-O6	-6.20	124.88	128.60
53	B5	499	G	C5-C6-O6	-6.20	124.88	128.60
53	B5	833	G	O4'-C1'-N9	6.20	113.16	108.20
1	AA	517	U	O4'-C1'-N1	6.20	113.16	108.20
53	B5	1089	G	O4'-C1'-N9	6.20	113.16	108.20
1	AA	597	G	N1-C6-O6	6.20	123.62	119.90
1	AA	816	G	O4'-C1'-N9	6.20	113.16	108.20
1	AA	1044	G	O4'-C1'-N9	6.20	113.16	108.20
53	B5	92	G	O4'-C1'-N9	6.20	113.16	108.20
53	B5	651	G	O4'-C1'-N9	6.20	113.16	108.20
1	AA	165	G	N1-C6-O6	6.19	123.61	119.90
1	AA	834	G	N1-C6-O6	6.19	123.61	119.90
53	B5	1542	G	O4'-C1'-N9	6.19	113.15	108.20
53	B5	2127	U	O4'-C1'-N1	6.19	113.15	108.20
1	AA	201	G	O4'-C1'-N9	6.19	113.15	108.20
1	AA	783	G	O4'-C1'-N9	6.19	113.15	108.20
1	AA	1126	U	O4'-C1'-N1	6.19	113.15	108.20
1	AA	1708	U	O4'-C1'-N1	6.19	113.15	108.20
53	B5	203	G	N1-C6-O6	6.19	123.61	119.90
53	B5	1006	A	O4'-C1'-N9	6.19	113.15	108.20
53	B5	1460	A	O4'-C1'-N9	6.19	113.15	108.20
53	B5	2671	A	O4'-C1'-N9	6.19	113.15	108.20
53	B5	2870	C	O4'-C1'-N1	6.19	113.15	108.20
1	AA	30	G	C5-C6-O6	-6.19	124.89	128.60
51	B3	85	G	O4'-C1'-N9	6.19	113.15	108.20
53	B5	1152	G	C5-C6-O6	-6.19	124.89	128.60
1	AA	600	U	O4'-C1'-N1	6.19	113.15	108.20
52	B4	62	C	C2-N1-C1'	6.19	125.61	118.80
53	B5	250	U	O4'-C1'-N1	6.19	113.15	108.20
53	B5	1266	G	O4'-C1'-N9	6.19	113.15	108.20
1	AA	858	G	C5'-C4'-O4'	6.19	116.52	109.10
1	AA	1214	G	N1-C6-O6	6.19	123.61	119.90
1	AA	701	U	O4'-C1'-N1	6.18	113.15	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1238	A	C5-C6-N6	-6.18	118.75	123.70
1	AA	1426	G	O4'-C1'-N9	6.18	113.15	108.20
53	B5	1166	G	O4'-C1'-N9	6.18	113.15	108.20
53	B5	3343	G	O4'-C1'-N9	6.18	113.15	108.20
1	AA	1394	U	O4'-C1'-N1	6.18	113.15	108.20
53	B5	1447	G	C5-C6-O6	-6.18	124.89	128.60
53	B5	2973	G	O4'-C1'-N9	6.18	113.15	108.20
1	AA	1296	G	C5-C6-O6	-6.18	124.89	128.60
51	B3	116	U	O4'-C1'-N1	6.18	113.14	108.20
53	B5	1808	G	O4'-C1'-N9	6.18	113.14	108.20
1	AA	753	A	O4'-C1'-N9	6.18	113.14	108.20
1	AA	1069	C	O4'-C1'-N1	6.18	113.14	108.20
1	AA	1260	G	O4'-C1'-N9	6.18	113.14	108.20
1	AA	1504	G	N1-C6-O6	6.18	123.61	119.90
19	A7	66	A	O4'-C4'-C3'	6.18	111.04	106.10
53	B5	371	G	O4'-C1'-N9	6.18	113.14	108.20
53	B5	1148	G	O4'-C1'-N9	6.18	113.14	108.20
53	B5	1944	U	O4'-C1'-N1	6.18	113.14	108.20
1	AA	1500	G	N1-C6-O6	6.18	123.61	119.90
53	B5	597	G	C5-C6-O6	-6.18	124.89	128.60
1	AA	528	U	O4'-C1'-N1	6.18	113.14	108.20
1	AA	1521	G	C6-N1-C2	6.18	128.81	125.10
1	AA	1775	G	C5-C6-O6	-6.18	124.89	128.60
52	B4	24	G	C5-C6-O6	-6.18	124.89	128.60
53	B5	1673	G	O4'-C1'-N9	6.18	113.14	108.20
1	AA	223	U	O4'-C1'-N1	6.17	113.14	108.20
1	AA	272	U	O4'-C1'-N1	6.17	113.14	108.20
1	AA	307	G	C5-C6-O6	-6.17	124.89	128.60
1	AA	1776	G	C5-C6-O6	-6.17	124.90	128.60
53	B5	128	G	N1-C6-O6	6.17	123.61	119.90
53	B5	963	G	C5-C6-O6	-6.17	124.90	128.60
53	B5	1392	G	C5-C6-O6	-6.17	124.89	128.60
1	AA	833	U	O4'-C1'-N1	6.17	113.14	108.20
1	AA	1508	U	O4'-C1'-N1	6.17	113.14	108.20
53	B5	1901	A	O4'-C1'-N9	6.17	113.14	108.20
1	AA	392	G	N1-C6-O6	6.17	123.60	119.90
53	B5	2370	G	C5-C6-O6	-6.17	124.90	128.60
19	A7	75	C	N3-C2-O2	-6.17	117.58	121.90
53	B5	1838	G	C5-C6-O6	-6.17	124.90	128.60
1	AA	1685	U	O4'-C1'-N1	6.17	113.13	108.20
1	AA	1696	G	O4'-C1'-N9	6.17	113.13	108.20
19	A7	9	A	C8-N9-C4	-6.17	103.33	105.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	251	G	C5-C6-O6	-6.17	124.90	128.60
53	B5	1543	G	C5-C6-O6	-6.17	124.90	128.60
53	B5	2954	U	O4'-C1'-N1	6.17	113.13	108.20
1	AA	216	U	O4'-C1'-N1	6.17	113.13	108.20
1	AA	370	A	O4'-C1'-N9	6.17	113.13	108.20
1	AA	957	U	O4'-C1'-N1	6.17	113.13	108.20
1	AA	975	G	O4'-C1'-N9	6.17	113.13	108.20
1	AA	1606	U	O4'-C1'-N1	6.17	113.13	108.20
1	AA	986	G	C5-C6-O6	-6.16	124.90	128.60
19	A7	56	C	O4'-C1'-C2'	6.16	113.15	107.60
48	BX	7	TYR	CB-CG-CD2	-6.16	117.30	121.00
52	B4	71	G	C5-C6-O6	-6.16	124.90	128.60
53	B5	74	G	C5-C6-O6	-6.16	124.90	128.60
53	B5	472	A	O4'-C1'-N9	6.16	113.13	108.20
53	B5	1090	G	O4'-C1'-N9	6.16	113.13	108.20
53	B5	2475	G	C5-C6-O6	-6.16	124.90	128.60
53	B5	725	G	N1-C6-O6	6.16	123.60	119.90
53	B5	2712	U	O4'-C1'-N1	6.16	113.13	108.20
53	B5	272	G	N1-C6-O6	6.16	123.60	119.90
53	B5	2503	G	O4'-C1'-N9	6.16	113.13	108.20
53	B5	2533	G	N1-C6-O6	6.16	123.60	119.90
1	AA	351	C	O4'-C1'-N1	6.16	113.13	108.20
1	AA	1281	C	O4'-C1'-N1	6.16	113.13	108.20
53	B5	717	C	C2-N1-C1'	6.16	125.58	118.80
53	B5	990	A	O4'-C1'-N9	6.16	113.13	108.20
53	B5	1929	G	N1-C6-O6	6.16	123.59	119.90
1	AA	898	G	C5-C6-O6	-6.16	124.91	128.60
1	AA	1677	G	N1-C6-O6	6.16	123.59	119.90
19	A7	74	C	N1-C2-O2	6.16	122.59	118.90
43	BS	21	PHE	CB-CG-CD2	6.16	125.11	120.80
53	B5	55	G	O4'-C1'-N9	6.16	113.12	108.20
53	B5	139	G	O4'-C1'-N9	6.16	113.12	108.20
53	B5	283	G	C5-C6-O6	-6.16	124.91	128.60
53	B5	2314	U	O4'-C1'-N1	6.16	113.12	108.20
53	B5	2973	G	C5-C6-O6	-6.16	124.91	128.60
53	B5	1695	U	O4'-C1'-N1	6.15	113.12	108.20
53	B5	2182	A	O4'-C1'-N9	6.15	113.12	108.20
53	B5	210	U	O4'-C1'-N1	6.15	113.12	108.20
53	B5	1795	U	O4'-C1'-N1	6.15	113.12	108.20
53	B5	1873	U	O4'-C1'-N1	6.15	113.12	108.20
53	B5	3251	U	O4'-C1'-N1	6.15	113.12	108.20
53	B5	3340	G	C5-C6-O6	-6.15	124.91	128.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1169	G	C5-C6-O6	-6.15	124.91	128.60
1	AA	1187	U	O4'-C1'-N1	6.15	113.12	108.20
1	AA	1667	U	O4'-C1'-N1	6.15	113.12	108.20
53	B5	989	U	O4'-C1'-N1	6.15	113.12	108.20
53	B5	1441	G	C5-C6-O6	-6.15	124.91	128.60
53	B5	1920	U	O4'-C1'-N1	6.15	113.12	108.20
53	B5	2660	G	O4'-C1'-N9	6.15	113.12	108.20
53	B5	3137	C	O4'-C1'-N1	6.15	113.12	108.20
53	B5	2824	G	N1-C6-O6	6.15	123.59	119.90
53	B5	3226	A	O4'-C1'-N9	6.15	113.12	108.20
1	AA	495	C	C6-N1-C1'	-6.15	113.42	120.80
53	B5	2822	U	O4'-C1'-N1	6.15	113.12	108.20
1	AA	177	U	O4'-C1'-N1	6.15	113.12	108.20
1	AA	595	G	C5-C6-O6	-6.15	124.91	128.60
53	B5	504	A	O4'-C1'-N9	6.15	113.12	108.20
53	B5	1878	G	C5-C6-O6	-6.15	124.91	128.60
53	B5	2874	G	O4'-C1'-N9	6.15	113.12	108.20
53	B5	3009	G	O4'-C1'-N9	6.15	113.12	108.20
1	AA	548	G	N1-C6-O6	6.14	123.59	119.90
1	AA	856	A	O3'-P-O5'	6.14	115.68	104.00
1	AA	1289	G	O4'-C1'-N9	6.14	113.11	108.20
53	B5	345	G	N1-C6-O6	6.14	123.59	119.90
53	B5	678	G	O4'-C1'-N9	6.14	113.11	108.20
53	B5	3353	G	O4'-C1'-N9	6.14	113.12	108.20
1	AA	32	U	O4'-C1'-N1	6.14	113.11	108.20
53	B5	2302	G	C5-C6-O6	-6.14	124.91	128.60
53	B5	3158	G	O4'-C1'-N9	6.14	113.11	108.20
1	AA	430	G	O4'-C1'-N9	6.14	113.11	108.20
53	B5	1573	G	C5-C6-O6	-6.14	124.92	128.60
53	B5	3073	A	O4'-C1'-N9	6.14	113.11	108.20
53	B5	452	G	C5-C6-O6	-6.14	124.92	128.60
53	B5	799	G	N1-C6-O6	6.14	123.58	119.90
53	B5	2174	G	C5-C6-O6	-6.14	124.92	128.60
53	B5	3085	G	C5-C6-O6	-6.14	124.92	128.60
53	B5	2268	U	O4'-C1'-N1	6.14	113.11	108.20
53	B5	2858	U	O4'-C1'-N1	6.14	113.11	108.20
53	B5	2900	A	O4'-C1'-N9	6.14	113.11	108.20
53	B5	530	G	O4'-C1'-N9	6.14	113.11	108.20
53	B5	822	G	O4'-C1'-N9	6.14	113.11	108.20
1	AA	419	G	O4'-C1'-N9	6.13	113.11	108.20
1	AA	522	U	O4'-C1'-N1	6.13	113.11	108.20
1	AA	586	G	O4'-C1'-N9	6.13	113.11	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	624	G	O4'-C1'-N9	6.13	113.11	108.20
53	B5	825	U	O4'-C1'-N1	6.13	113.11	108.20
1	AA	15	U	O4'-C1'-N1	6.13	113.11	108.20
1	AA	25	C	O4'-C1'-N1	6.13	113.11	108.20
1	AA	516	G	O4'-C1'-N9	6.13	113.11	108.20
53	B5	2781	U	O4'-C1'-N1	6.13	113.11	108.20
1	AA	109	G	N1-C6-O6	6.13	123.58	119.90
1	AA	113	U	O4'-C1'-N1	6.13	113.10	108.20
1	AA	921	G	O4'-C1'-N9	6.13	113.10	108.20
1	AA	993	G	N1-C6-O6	6.13	123.58	119.90
1	AA	1452	G	O4'-C1'-N9	6.13	113.10	108.20
53	B5	216	G	O4'-C1'-N9	6.13	113.10	108.20
53	B5	685	G	N1-C6-O6	6.13	123.58	119.90
53	B5	1004	U	O4'-C1'-N1	6.13	113.10	108.20
53	B5	2400	G	C5-C6-O6	-6.13	124.92	128.60
1	AA	1268	U	O4'-C1'-N1	6.13	113.10	108.20
1	AA	1722	U	O4'-C1'-N1	6.13	113.10	108.20
53	B5	848	A	O4'-C1'-N9	6.13	113.10	108.20
53	B5	1206	G	O4'-C1'-N9	6.13	113.10	108.20
53	B5	2677	G	C5-C6-O6	-6.13	124.92	128.60
1	AA	1163	A	O4'-C1'-N9	6.12	113.10	108.20
53	B5	1547	G	N1-C6-O6	6.12	123.58	119.90
53	B5	2246	G	N1-C6-O6	6.12	123.58	119.90
53	B5	2573	G	N1-C6-O6	6.12	123.58	119.90
53	B5	3392	U	O4'-C1'-N1	6.12	113.10	108.20
1	AA	313	U	O4'-C1'-N1	6.12	113.10	108.20
1	AA	640	U	O4'-C1'-N1	6.12	113.10	108.20
1	AA	1311	U	O4'-C1'-N1	6.12	113.10	108.20
53	B5	35	A	O4'-C1'-N9	6.12	113.10	108.20
53	B5	267	G	N1-C6-O6	6.12	123.57	119.90
53	B5	1691	U	O4'-C1'-N1	6.12	113.10	108.20
53	B5	2605	G	C5-C6-O6	-6.12	124.93	128.60
1	AA	647	G	O4'-C1'-N9	6.12	113.10	108.20
1	AA	813	U	O4'-C1'-N1	6.12	113.10	108.20
1	AA	1002	G	O4'-C1'-N9	6.12	113.10	108.20
1	AA	1341	A	C1'-O4'-C4'	-6.12	105.00	109.90
1	AA	1521	G	C5-N7-C8	-6.12	101.24	104.30
1	AA	1551	G	O4'-C4'-C3'	-6.12	97.88	104.00
19	A7	48	C	N1-C2-O2	6.12	122.57	118.90
53	B5	1024	G	C5-C6-O6	-6.12	124.93	128.60
53	B5	2633	U	O4'-C1'-N1	6.12	113.10	108.20
53	B5	2660	G	N1-C6-O6	6.12	123.57	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	514	G	O4'-C1'-N9	6.12	113.09	108.20
1	AA	514	G	N1-C6-O6	6.12	123.57	119.90
1	AA	1031	G	N1-C6-O6	6.12	123.57	119.90
1	AA	1108	G	N1-C6-O6	6.12	123.57	119.90
1	AA	1553	A	O5'-P-OP2	6.12	118.04	110.70
1	AA	1756	U	O4'-C1'-N1	6.12	113.09	108.20
52	B4	116	G	C5-C6-O6	-6.12	124.93	128.60
53	B5	1382	G	O4'-C1'-N9	6.12	113.09	108.20
53	B5	3191	G	O4'-C1'-N9	6.12	113.09	108.20
1	AA	1043	G	C5-C6-O6	-6.12	124.93	128.60
53	B5	3209	U	O4'-C1'-N1	6.12	113.09	108.20
1	AA	845	G	N1-C6-O6	6.12	123.57	119.90
19	A7	75	C	C4'-C3'-C2'	-6.12	96.48	102.60
53	B5	2246	G	O4'-C1'-N9	6.12	113.09	108.20
1	AA	1367	U	O4'-C1'-N1	6.11	113.09	108.20
19	A7	74	C	O4'-C4'-C3'	6.11	110.99	106.10
53	B5	227	G	O4'-C1'-N9	6.11	113.09	108.20
53	B5	2687	G	C5-C6-O6	-6.11	124.93	128.60
53	B5	3091	A	O4'-C1'-N9	6.11	113.09	108.20
53	B5	3318	G	O4'-C1'-N9	6.11	113.09	108.20
1	AA	807	A	O4'-C1'-N9	6.11	113.09	108.20
53	B5	853	G	N1-C6-O6	6.11	123.57	119.90
1	AA	563	U	O4'-C1'-N1	6.11	113.09	108.20
1	AA	1523	A	N1-C6-N6	6.11	122.27	118.60
1	AA	1551	G	C2'-C3'-O3'	-6.11	96.06	109.50
53	B5	353	G	O4'-C1'-N9	6.11	113.09	108.20
53	B5	958	C	O4'-C1'-N1	6.11	113.09	108.20
53	B5	2922	G	O4'-C1'-N9	6.11	113.09	108.20
53	B5	3269	A	O4'-C1'-N9	6.11	113.09	108.20
53	B5	2369	G	O4'-C1'-N9	6.11	113.09	108.20
53	B5	197	G	O4'-C1'-N9	6.11	113.09	108.20
53	B5	2104	A	O4'-C1'-N9	6.11	113.09	108.20
53	B5	2922	G	C5-C6-O6	-6.11	124.94	128.60
53	B5	3177	U	O4'-C1'-N1	6.11	113.09	108.20
1	AA	860	U	C2'-C3'-O3'	-6.11	96.07	109.50
53	B5	530	G	N1-C6-O6	6.11	123.56	119.90
53	B5	568	G	C5-C6-O6	-6.11	124.94	128.60
53	B5	964	G	N1-C6-O6	6.11	123.56	119.90
53	B5	1222	G	O4'-C1'-N9	6.11	113.08	108.20
53	B5	2669	G	N1-C6-O6	6.11	123.56	119.90
53	B5	2938	G	N1-C6-O6	6.11	123.56	119.90
1	AA	859	A	C4-C5-C6	6.10	120.05	117.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	866	G	O4'-C4'-C3'	-6.10	97.90	104.00
53	B5	419	G	O4'-C1'-N9	6.10	113.08	108.20
53	B5	552	G	O4'-C1'-N9	6.10	113.08	108.20
53	B5	684	G	N1-C6-O6	6.10	123.56	119.90
53	B5	820	A	C5-C6-N6	-6.10	118.82	123.70
53	B5	2848	G	O4'-C1'-N9	6.10	113.08	108.20
53	B5	3260	G	N1-C6-O6	6.10	123.56	119.90
1	AA	140	A	O4'-C1'-N9	6.10	113.08	108.20
1	AA	450	U	O4'-C1'-N1	6.10	113.08	108.20
53	B5	2107	A	O4'-C1'-N9	6.10	113.08	108.20
53	B5	2433	U	O4'-C1'-N1	6.10	113.08	108.20
53	B5	2714	G	C5-C6-O6	-6.10	124.94	128.60
53	B5	510	G	O4'-C1'-N9	6.10	113.08	108.20
53	B5	2794	G	C5-C6-O6	-6.10	124.94	128.60
1	AA	74	U	O4'-C1'-N1	6.10	113.08	108.20
52	B4	135	G	C5-C6-O6	-6.10	124.94	128.60
53	B5	583	G	C5-C6-O6	-6.10	124.94	128.60
53	B5	986	U	O4'-C1'-N1	6.10	113.08	108.20
53	B5	1933	A	C5-C6-N6	-6.10	118.82	123.70
1	AA	429	G	N1-C6-O6	6.10	123.56	119.90
1	AA	981	U	O4'-C1'-N1	6.10	113.08	108.20
53	B5	604	G	C5-C6-O6	-6.10	124.94	128.60
53	B5	1598	G	O4'-C1'-N9	6.10	113.08	108.20
53	B5	1903	U	O4'-C1'-N1	6.10	113.08	108.20
53	B5	2608	G	O4'-C1'-N9	6.10	113.08	108.20
53	B5	1019	G	N1-C6-O6	6.10	123.56	119.90
1	AA	199	G	O4'-C1'-N9	6.09	113.08	108.20
52	B4	56	G	O4'-C1'-N9	6.09	113.08	108.20
53	B5	511	G	O4'-C1'-N9	6.09	113.07	108.20
53	B5	1413	G	C5-C6-O6	-6.09	124.94	128.60
53	B5	3044	G	N1-C6-O6	6.09	123.56	119.90
1	AA	597	G	O4'-C1'-N9	6.09	113.07	108.20
19	A7	11	C	C2-N3-C4	-6.09	116.85	119.90
53	B5	474	G	O4'-C1'-N9	6.09	113.07	108.20
53	B5	1178	G	C5-C6-O6	-6.09	124.94	128.60
53	B5	2713	U	O4'-C1'-N1	6.09	113.07	108.20
1	AA	331	A	O4'-C1'-N9	6.09	113.07	108.20
1	AA	761	G	C5-C6-O6	-6.09	124.95	128.60
1	AA	1251	G	N1-C6-O6	6.09	123.55	119.90
1	AA	1502	G	N1-C6-O6	6.09	123.55	119.90
1	AA	1661	G	N1-C6-O6	6.09	123.55	119.90
52	B4	95	G	N1-C6-O6	6.09	123.55	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	1471	U	O4'-C1'-N1	6.09	113.07	108.20
53	B5	3205	U	O4'-C1'-N1	6.09	113.07	108.20
1	AA	762	A	O4'-C1'-N9	6.09	113.07	108.20
53	B5	3068	U	O4'-C1'-N1	6.09	113.07	108.20
1	AA	544	A	O4'-C1'-N9	6.09	113.07	108.20
1	AA	1584	A	C4-C5-C6	6.09	120.04	117.00
1	AA	1585	A	C5-C6-N6	-6.09	118.83	123.70
1	AA	1590	A	O4'-C1'-N9	6.09	113.07	108.20
1	AA	1694	G	N1-C6-O6	6.09	123.55	119.90
51	B3	64	A	O4'-C1'-N9	6.09	113.07	108.20
53	B5	1235	U	O4'-C1'-N1	6.09	113.07	108.20
53	B5	1400	G	O4'-C1'-N9	6.09	113.07	108.20
53	B5	1652	G	C5-C6-O6	-6.09	124.95	128.60
53	B5	2970	C	O4'-C1'-N1	6.09	113.07	108.20
53	B5	2997	G	O4'-C1'-N9	6.08	113.07	108.20
1	AA	1372	A	O4'-C1'-N9	6.08	113.07	108.20
1	AA	1635	C	C5'-C4'-O4'	-6.08	101.80	109.10
19	A7	62	A	C1'-O4'-C4'	-6.08	105.03	109.90
53	B5	229	G	N1-C6-O6	6.08	123.55	119.90
1	AA	75	U	O4'-C1'-N1	6.08	113.07	108.20
1	AA	576	G	N1-C6-O6	6.08	123.55	119.90
1	AA	592	A	O4'-C1'-N9	6.08	113.06	108.20
53	B5	1227	C	P-O3'-C3'	6.08	127.00	119.70
53	B5	2452	G	C5-C6-O6	-6.08	124.95	128.60
1	AA	744	U	O4'-C1'-N1	6.08	113.06	108.20
1	AA	1745	G	O4'-C1'-N9	6.08	113.06	108.20
53	B5	2315	G	O4'-C1'-N9	6.08	113.06	108.20
53	B5	3287	U	O4'-C1'-N1	6.08	113.06	108.20
1	AA	371	G	O4'-C1'-N9	6.08	113.06	108.20
1	AA	1724	G	O4'-C1'-N9	6.08	113.06	108.20
53	B5	2441	A	O4'-C1'-N9	6.08	113.06	108.20
1	AA	337	G	N1-C6-O6	6.08	123.55	119.90
53	B5	647	A	O4'-C1'-N9	6.08	113.06	108.20
53	B5	1341	U	O4'-C1'-N1	6.08	113.06	108.20
1	AA	608	U	O4'-C1'-N1	6.07	113.06	108.20
1	AA	1235	U	O4'-C1'-N1	6.07	113.06	108.20
53	B5	1241	U	O4'-C1'-N1	6.07	113.06	108.20
53	B5	1804	A	O4'-C1'-N9	6.07	113.06	108.20
53	B5	2193	U	O4'-C1'-N1	6.07	113.06	108.20
28	BD	194	TYR	CB-CG-CD1	-6.07	117.36	121.00
53	B5	770	G	O4'-C1'-N9	6.07	113.06	108.20
53	B5	1599	G	O4'-C1'-N9	6.07	113.06	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	1655	G	C5-C6-O6	-6.07	124.96	128.60
1	AA	1377	U	O4'-C1'-N1	6.07	113.06	108.20
53	B5	2936	A	O4'-C1'-N9	6.07	113.06	108.20
53	B5	1262	G	N1-C6-O6	6.07	123.54	119.90
1	AA	810	G	N1-C6-O6	6.07	123.54	119.90
1	AA	1047	G	N1-C6-O6	6.07	123.54	119.90
1	AA	1460	G	O4'-C1'-N9	6.07	113.05	108.20
1	AA	1735	G	O4'-C1'-N9	6.07	113.05	108.20
50	BZ	43	TYR	CB-CG-CD2	-6.07	117.36	121.00
51	B3	107	G	C5-C6-O6	-6.07	124.96	128.60
53	B5	11	A	O4'-C1'-N9	6.07	113.05	108.20
53	B5	778	U	O4'-C1'-N1	6.07	113.06	108.20
1	AA	497	G	O4'-C1'-N9	6.07	113.05	108.20
1	AA	541	A	O4'-C1'-N9	6.07	113.05	108.20
1	AA	655	G	O4'-C1'-N9	6.06	113.05	108.20
19	A7	66	A	N1-C2-N3	-6.06	126.27	129.30
53	B5	1650	G	O4'-C1'-N9	6.06	113.05	108.20
1	AA	557	G	N1-C6-O6	6.06	123.54	119.90
1	AA	881	A	O4'-C1'-N9	6.06	113.05	108.20
38	BN	60	PHE	CB-CG-CD1	6.06	125.04	120.80
53	B5	805	G	C5-C6-O6	-6.06	124.96	128.60
53	B5	908	G	C5-C6-O6	-6.06	124.96	128.60
53	B5	1278	A	O4'-C1'-N9	6.06	113.05	108.20
53	B5	197	G	C5-C6-O6	-6.06	124.96	128.60
1	AA	1106	G	N1-C6-O6	6.06	123.54	119.90
52	B4	36	G	C5-C6-O6	-6.06	124.97	128.60
53	B5	468	G	N1-C6-O6	6.06	123.53	119.90
53	B5	770	G	C5-C6-O6	-6.06	124.96	128.60
53	B5	2121	G	O4'-C1'-N9	6.06	113.05	108.20
53	B5	2553	U	O4'-C1'-N1	6.06	113.05	108.20
53	B5	3199	G	O4'-C1'-N9	6.06	113.05	108.20
1	AA	9	U	O4'-C1'-N1	6.06	113.05	108.20
1	AA	290	G	N1-C6-O6	6.06	123.53	119.90
1	AA	466	U	O4'-C1'-N1	6.06	113.05	108.20
1	AA	739	G	O4'-C1'-N9	6.06	113.05	108.20
1	AA	918	A	O4'-C1'-N9	6.06	113.05	108.20
1	AA	923	A	O4'-C1'-N9	6.06	113.05	108.20
1	AA	1688	G	O4'-C1'-N9	6.06	113.05	108.20
53	B5	960	U	O4'-C1'-N1	6.06	113.05	108.20
53	B5	1288	U	O4'-C1'-N1	6.06	113.05	108.20
53	B5	2425	G	C5-C6-O6	-6.06	124.97	128.60
1	AA	332	U	O4'-C1'-N1	6.06	113.05	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	1253	U	O4'-C1'-N1	6.06	113.05	108.20
53	B5	1897	G	O4'-C1'-N9	6.06	113.05	108.20
51	B3	49	G	O4'-C1'-N9	6.05	113.04	108.20
53	B5	244	G	O4'-C1'-N9	6.05	113.04	108.20
1	AA	336	G	O4'-C1'-N9	6.05	113.04	108.20
53	B5	1177	G	C5-C6-O6	-6.05	124.97	128.60
53	B5	1182	A	O4'-C1'-N9	6.05	113.04	108.20
53	B5	3131	U	O4'-C1'-N1	6.05	113.04	108.20
52	B4	132	G	O4'-C1'-N9	6.05	113.04	108.20
53	B5	1144	U	O4'-C1'-N1	6.05	113.04	108.20
1	AA	319	U	C2-N1-C1'	6.05	124.96	117.70
43	BS	21	PHE	CB-CG-CD1	-6.05	116.56	120.80
53	B5	1409	G	O4'-C1'-N9	6.05	113.04	108.20
1	AA	115	G	C5-C6-O6	-6.05	124.97	128.60
53	B5	445	G	O4'-C1'-N9	6.05	113.04	108.20
53	B5	1813	A	O4'-C1'-N9	6.05	113.04	108.20
53	B5	2535	A	O4'-C1'-N9	6.05	113.04	108.20
53	B5	998	A	O4'-C1'-N9	6.05	113.04	108.20
53	B5	2860	U	O4'-C1'-N1	6.05	113.04	108.20
1	AA	1108	G	O4'-C1'-N9	6.04	113.04	108.20
1	AA	808	U	O4'-C1'-N1	6.04	113.03	108.20
1	AA	998	U	O4'-C1'-N1	6.04	113.03	108.20
1	AA	1188	C	O4'-C1'-N1	6.04	113.03	108.20
53	B5	1021	G	O4'-C1'-N9	6.04	113.03	108.20
53	B5	2180	G	C5-C6-O6	-6.04	124.97	128.60
53	B5	2377	G	N1-C6-O6	6.04	123.53	119.90
53	B5	2573	G	O4'-C1'-N9	6.04	113.03	108.20
53	B5	2867	C	O4'-C1'-N1	6.04	113.03	108.20
1	AA	161	U	O4'-C1'-N1	6.04	113.03	108.20
1	AA	844	A	O4'-C1'-N9	6.04	113.03	108.20
19	A7	56	C	N1-C2-N3	-6.04	114.97	119.20
19	A7	61	C	C5'-C4'-O4'	6.04	116.35	109.10
1	AA	899	A	P-O3'-C3'	6.04	126.95	119.70
53	B5	514	G	O4'-C1'-N9	6.04	113.03	108.20
53	B5	2680	A	O4'-C1'-N9	6.04	113.03	108.20
1	AA	10	G	N1-C6-O6	6.04	123.52	119.90
1	AA	973	A	O4'-C1'-N9	6.04	113.03	108.20
1	AA	1655	U	O4'-C1'-N1	6.04	113.03	108.20
19	A7	57	G	N1-C2-N3	6.04	127.52	123.90
51	B3	51	G	C5-C6-O6	-6.04	124.98	128.60
52	B4	97	A	O4'-C1'-N9	6.04	113.03	108.20
53	B5	737	G	C5-C6-O6	-6.04	124.98	128.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	1188	U	O4'-C1'-N1	6.04	113.03	108.20
53	B5	1934	G	O4'-C1'-N9	6.04	113.03	108.20
53	B5	3381	U	O4'-C1'-N1	6.04	113.03	108.20
1	AA	1453	G	O4'-C1'-N9	6.04	113.03	108.20
53	B5	864	G	C5-C6-O6	-6.04	124.98	128.60
53	B5	2278	C	C6-N1-C2	-6.04	117.89	120.30
53	B5	2390	A	O4'-C1'-N9	6.04	113.03	108.20
53	B5	2746	A	O4'-C1'-N9	6.04	113.03	108.20
53	B5	3075	G	C5-C6-O6	-6.04	124.98	128.60
53	B5	596	U	O4'-C1'-N1	6.03	113.03	108.20
19	A7	11	C	N3-C4-N4	-6.03	113.78	118.00
53	B5	1675	G	C5-C6-O6	-6.03	124.98	128.60
53	B5	1796	G	C5-C6-O6	-6.03	124.98	128.60
1	AA	395	U	O4'-C1'-N1	6.03	113.02	108.20
1	AA	1659	U	O4'-C1'-N1	6.03	113.03	108.20
53	B5	920	A	C5-C6-N6	-6.03	118.88	123.70
53	B5	2732	G	C5-C6-O6	-6.03	124.98	128.60
1	AA	879	G	N1-C6-O6	6.03	123.52	119.90
53	B5	252	U	O4'-C1'-N1	6.03	113.02	108.20
53	B5	595	A	O4'-C1'-N9	6.03	113.02	108.20
53	B5	1166	G	N1-C6-O6	6.03	123.52	119.90
53	B5	2796	G	O4'-C1'-N9	6.03	113.02	108.20
53	B5	629	U	O4'-C1'-N1	6.03	113.02	108.20
1	AA	1414	A	O4'-C1'-N9	6.02	113.02	108.20
53	B5	3266	G	O4'-C1'-N9	6.02	113.02	108.20
1	AA	801	G	O4'-C1'-N9	6.02	113.02	108.20
1	AA	1713	G	O4'-C1'-N9	6.02	113.02	108.20
53	B5	3395	G	N1-C6-O6	6.02	123.51	119.90
1	AA	221	A	O4'-C1'-N9	6.02	113.02	108.20
1	AA	676	G	O4'-C1'-N9	6.02	113.02	108.20
51	B3	88	G	O4'-C1'-N9	6.02	113.02	108.20
53	B5	915	A	O4'-C1'-N9	6.02	113.02	108.20
1	AA	1132	U	O4'-C1'-N1	6.02	113.02	108.20
1	AA	1788	A	C5-C6-N6	-6.02	118.88	123.70
53	B5	267	G	O4'-C1'-N9	6.02	113.02	108.20
53	B5	1534	A	O4'-C1'-N9	6.02	113.02	108.20
53	B5	1613	A	O4'-C1'-N9	6.02	113.02	108.20
53	B5	2470	C	C6-N1-C1'	-6.02	113.58	120.80
53	B5	2996	U	O4'-C1'-N1	6.02	113.02	108.20
1	AA	994	A	O4'-C1'-N9	6.02	113.02	108.20
1	AA	1748	A	O4'-C1'-N9	6.02	113.01	108.20
53	B5	1546	A	O4'-C1'-N9	6.02	113.01	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	1661	G	O4'-C1'-N9	6.02	113.01	108.20
53	B5	2961	G	C5-C6-O6	-6.02	124.99	128.60
1	AA	497	G	N1-C6-O6	6.02	123.51	119.90
1	AA	44	U	O4'-C1'-N1	6.01	113.01	108.20
1	AA	1634	C	O4'-C1'-N1	6.01	113.01	108.20
1	AA	1711	G	O4'-C1'-N9	6.01	113.01	108.20
52	B4	92	A	C5-C6-N6	-6.01	118.89	123.70
53	B5	483	G	N1-C6-O6	6.01	123.51	119.90
53	B5	905	U	O4'-C1'-N1	6.01	113.01	108.20
53	B5	1736	G	C5-C6-O6	-6.01	124.99	128.60
53	B5	1823	A	P-O3'-C3'	6.01	126.92	119.70
53	B5	2576	G	O4'-C1'-N9	6.01	113.01	108.20
1	AA	512	A	O4'-C1'-N9	6.01	113.01	108.20
1	AA	1746	G	O4'-C1'-N9	6.01	113.01	108.20
53	B5	2275	A	O4'-C1'-N9	6.01	113.01	108.20
1	AA	670	U	O4'-C1'-N1	6.01	113.01	108.20
19	A7	41	U	N1-C2-O2	-6.01	118.59	122.80
53	B5	810	A	C5-C6-N6	-6.01	118.89	123.70
53	B5	1148	G	N1-C6-O6	6.01	123.51	119.90
53	B5	2739	A	O4'-C1'-N9	6.01	113.01	108.20
53	B5	3160	U	O4'-C1'-N1	6.01	113.01	108.20
53	B5	3281	U	O4'-C1'-N1	6.01	113.01	108.20
1	AA	1475	G	N1-C6-O6	6.01	123.50	119.90
1	AA	1225	G	O4'-C1'-N9	6.01	113.01	108.20
1	AA	1773	U	O4'-C1'-N1	6.01	113.00	108.20
53	B5	1012	G	C5-C6-O6	-6.01	125.00	128.60
53	B5	2361	A	O4'-C1'-N9	6.01	113.00	108.20
1	AA	1530	U	O4'-C1'-N1	6.00	113.00	108.20
19	A7	27	C	O4'-C1'-N1	6.00	113.00	108.20
53	B5	867	G	C5-C6-O6	-6.00	125.00	128.60
53	B5	1124	U	O4'-C1'-N1	6.00	113.00	108.20
1	AA	349	U	O4'-C1'-N1	6.00	113.00	108.20
1	AA	1515	U	O3'-P-O5'	-6.00	92.60	104.00
52	B4	73	U	O4'-C1'-N1	6.00	113.00	108.20
53	B5	1387	G	O4'-C1'-N9	6.00	113.00	108.20
53	B5	2158	A	O4'-C1'-N9	6.00	113.00	108.20
53	B5	2589	G	C5-C6-O6	-6.00	125.00	128.60
53	B5	300	G	O4'-C1'-N9	6.00	113.00	108.20
1	AA	1359	U	O4'-C1'-N1	6.00	113.00	108.20
53	B5	1133	A	O4'-C1'-N9	6.00	113.00	108.20
53	B5	1447	G	P-O3'-C3'	6.00	126.90	119.70
1	AA	233	C	O4'-C1'-N1	6.00	113.00	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	398	G	C5-C6-O6	-6.00	125.00	128.60
1	AA	749	U	O4'-C1'-N1	6.00	113.00	108.20
1	AA	1614	G	O4'-C1'-N9	6.00	113.00	108.20
53	B5	459	G	O4'-C1'-N9	6.00	113.00	108.20
1	AA	1516	C	O4'-C4'-C3'	-6.00	98.00	104.00
1	AA	617	U	O4'-C1'-N1	5.99	113.00	108.20
1	AA	725	U	O4'-C1'-N1	5.99	112.99	108.20
1	AA	1099	G	C5-C6-O6	-5.99	125.00	128.60
53	B5	870	G	N1-C6-O6	5.99	123.50	119.90
53	B5	1020	G	O4'-C1'-N9	5.99	113.00	108.20
53	B5	2276	G	O4'-C1'-N9	5.99	113.00	108.20
53	B5	2429	G	O4'-C1'-N9	5.99	113.00	108.20
1	AA	388	G	N1-C6-O6	5.99	123.50	119.90
1	AA	599	A	O4'-C1'-N9	5.99	112.99	108.20
53	B5	527	A	O4'-C1'-N9	5.99	112.99	108.20
53	B5	2955	U	O4'-C1'-N1	5.99	112.99	108.20
1	AA	198	A	O4'-C1'-N9	5.99	112.99	108.20
1	AA	1366	U	O4'-C1'-N1	5.99	112.99	108.20
53	B5	381	U	O4'-C1'-N1	5.99	112.99	108.20
53	B5	1018	G	O4'-C1'-N9	5.99	112.99	108.20
1	AA	355	G	O4'-C1'-N9	5.99	112.99	108.20
1	AA	738	G	O4'-C1'-N9	5.99	112.99	108.20
1	AA	757	A	O4'-C1'-N9	5.99	112.99	108.20
53	B5	812	G	N1-C6-O6	5.99	123.49	119.90
53	B5	1040	A	O4'-C1'-N9	5.99	112.99	108.20
53	B5	2856	G	O4'-C1'-N9	5.99	112.99	108.20
1	AA	117	U	O4'-C1'-N1	5.99	112.99	108.20
1	AA	406	U	O4'-C1'-N1	5.99	112.99	108.20
53	B5	316	U	O4'-C1'-N1	5.99	112.99	108.20
53	B5	360	G	O4'-C1'-N9	5.99	112.99	108.20
53	B5	1456	A	C4-C5-C6	5.99	119.99	117.00
53	B5	2950	G	N1-C6-O6	5.99	123.49	119.90
53	B5	947	G	O4'-C1'-N9	5.99	112.99	108.20
53	B5	1440	G	O4'-C1'-N9	5.99	112.99	108.20
1	AA	472	U	O4'-C1'-N1	5.98	112.99	108.20
53	B5	358	G	O4'-C1'-N9	5.98	112.99	108.20
1	AA	330	G	O4'-C1'-N9	5.98	112.99	108.20
1	AA	372	G	C5-C6-O6	-5.98	125.01	128.60
1	AA	356	G	O4'-C1'-N9	5.98	112.98	108.20
1	AA	476	U	O4'-C1'-N1	5.98	112.98	108.20
51	B3	98	G	C5-C6-O6	-5.98	125.01	128.60
53	B5	1599	G	N1-C6-O6	5.98	123.49	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	1640	G	N1-C6-O6	5.98	123.49	119.90
53	B5	3258	U	P-O3'-C3'	5.98	126.88	119.70
1	AA	1141	U	O4'-C1'-N1	5.98	112.98	108.20
1	AA	1787	G	C5-C6-O6	-5.98	125.01	128.60
53	B5	1519	G	C5-C6-O6	-5.98	125.01	128.60
1	AA	118	U	O4'-C1'-N1	5.98	112.98	108.20
1	AA	537	G	O4'-C1'-N9	5.98	112.98	108.20
1	AA	767	U	O4'-C1'-N1	5.98	112.98	108.20
53	B5	216	G	N1-C6-O6	5.98	123.49	119.90
53	B5	721	G	O4'-C1'-N9	5.98	112.98	108.20
53	B5	3328	G	O4'-C1'-N9	5.98	112.98	108.20
1	AA	1018	A	O4'-C1'-N9	5.98	112.98	108.20
52	B4	101	U	O4'-C1'-N1	5.98	112.98	108.20
53	B5	1209	G	C5-C6-O6	-5.98	125.01	128.60
1	AA	7	G	C5-C6-O6	-5.97	125.02	128.60
1	AA	299	A	O4'-C1'-N9	5.97	112.98	108.20
1	AA	775	G	O4'-C1'-N9	5.97	112.98	108.20
1	AA	235	G	C5-C6-O6	-5.97	125.02	128.60
53	B5	1020	G	N1-C6-O6	5.97	123.48	119.90
53	B5	2278	C	O4'-C1'-N1	5.97	112.98	108.20
53	B5	2603	G	N1-C6-O6	5.97	123.48	119.90
53	B5	3224	G	N1-C6-O6	5.97	123.48	119.90
53	B5	30	G	O4'-C1'-N9	5.97	112.98	108.20
1	AA	280	U	O4'-C1'-N1	5.97	112.98	108.20
1	AA	656	G	O4'-C1'-N9	5.97	112.98	108.20
1	AA	772	G	N1-C6-O6	5.97	123.48	119.90
1	AA	983	G	N1-C6-O6	5.97	123.48	119.90
53	B5	745	C	N3-C4-N4	5.97	122.18	118.00
53	B5	1598	G	N1-C6-O6	5.97	123.48	119.90
53	B5	1818	U	O4'-C1'-N1	5.97	112.98	108.20
53	B5	1892	G	O4'-C1'-N9	5.97	112.97	108.20
53	B5	2376	G	O4'-C1'-N9	5.97	112.98	108.20
53	B5	2670	G	O4'-C1'-N9	5.97	112.98	108.20
1	AA	748	U	O4'-C1'-N1	5.97	112.97	108.20
53	B5	2602	G	O4'-C1'-N9	5.97	112.97	108.20
53	B5	3393	U	O4'-C1'-N1	5.97	112.97	108.20
53	B5	1	G	O4'-C1'-N9	5.97	112.97	108.20
53	B5	165	A	O4'-C1'-N9	5.97	112.97	108.20
53	B5	401	U	O4'-C1'-N1	5.97	112.97	108.20
53	B5	2918	G	N1-C6-O6	5.97	123.48	119.90
1	AA	356	G	N1-C6-O6	5.96	123.48	119.90
1	AA	712	G	O4'-C1'-N9	5.96	112.97	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	850	A	O4'-C1'-N9	5.96	112.97	108.20
1	AA	1273	G	C5-C6-O6	-5.96	125.02	128.60
51	B3	7	G	N1-C6-O6	5.96	123.48	119.90
51	B3	25	G	O4'-C1'-N9	5.96	112.97	108.20
53	B5	393	U	O4'-C1'-N1	5.96	112.97	108.20
53	B5	999	G	O4'-C1'-N9	5.96	112.97	108.20
53	B5	1750	A	O4'-C1'-N9	5.96	112.97	108.20
53	B5	2602	G	N1-C6-O6	5.96	123.48	119.90
53	B5	2606	G	N1-C6-O6	5.96	123.48	119.90
53	B5	3009	G	C5-C6-O6	-5.96	125.02	128.60
1	AA	577	G	O4'-C1'-N9	5.96	112.97	108.20
1	AA	728	U	O4'-C1'-N1	5.96	112.97	108.20
1	AA	730	G	C5-C6-O6	-5.96	125.02	128.60
53	B5	415	G	C5-C6-O6	-5.96	125.02	128.60
53	B5	726	G	O4'-C1'-N9	5.96	112.97	108.20
1	AA	593	U	O4'-C1'-N1	5.96	112.97	108.20
1	AA	1031	G	O4'-C1'-N9	5.96	112.97	108.20
53	B5	470	G	N1-C6-O6	5.96	123.48	119.90
53	B5	1242	G	O4'-C1'-N9	5.96	112.97	108.20
53	B5	2898	G	C5-C6-O6	-5.96	125.02	128.60
53	B5	3036	G	N1-C6-O6	5.96	123.48	119.90
53	B5	518	G	O4'-C1'-N9	5.96	112.97	108.20
1	AA	1522	A	O5'-P-OP2	-5.96	100.34	105.70
53	B5	2339	C	O4'-C1'-N1	5.96	112.97	108.20
53	B5	2536	A	O4'-C1'-N9	5.96	112.97	108.20
53	B5	2877	G	C5-C6-O6	-5.96	125.03	128.60
1	AA	98	U	O4'-C1'-N1	5.96	112.97	108.20
1	AA	1200	A	C5-N7-C8	5.96	106.88	103.90
1	AA	1219	A	O4'-C1'-N9	5.96	112.97	108.20
53	B5	575	G	O4'-C1'-N9	5.96	112.97	108.20
53	B5	1906	G	N1-C6-O6	5.96	123.47	119.90
53	B5	2238	G	C5-C6-O6	-5.96	125.03	128.60
53	B5	3315	G	C5-C6-O6	-5.96	125.03	128.60
53	B5	3353	G	N1-C6-O6	5.96	123.47	119.90
1	AA	790	U	O4'-C1'-N1	5.96	112.96	108.20
53	B5	392	G	C5-C6-O6	-5.96	125.03	128.60
53	B5	769	G	O4'-C1'-N9	5.96	112.96	108.20
53	B5	858	A	O4'-C1'-N9	5.96	112.96	108.20
53	B5	2947	G	O4'-C1'-N9	5.95	112.96	108.20
53	B5	1496	C	O4'-C1'-N1	5.95	112.96	108.20
53	B5	773	G	C5-C6-O6	-5.95	125.03	128.60
53	B5	1306	G	N1-C6-O6	5.95	123.47	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	2206	G	C5-C6-O6	-5.95	125.03	128.60
53	B5	2751	G	C5-C6-O6	-5.95	125.03	128.60
53	B5	3178	U	O4'-C1'-N1	5.95	112.96	108.20
1	AA	1	U	C2-N1-C1'	5.95	124.84	117.70
1	AA	377	G	C5-C6-O6	-5.95	125.03	128.60
1	AA	503	G	C5-C6-O6	-5.95	125.03	128.60
1	AA	887	A	O4'-C1'-N9	5.95	112.96	108.20
1	AA	1411	U	O4'-C1'-N1	5.95	112.96	108.20
19	A7	9	A	C2-N3-C4	5.95	113.57	110.60
19	A7	56	C	C1'-O4'-C4'	-5.95	105.14	109.90
53	B5	619	A	C5-C6-N6	-5.95	118.94	123.70
53	B5	860	G	C4-N9-C1'	5.95	134.23	126.50
53	B5	1357	G	C5-C6-O6	-5.95	125.03	128.60
53	B5	2135	U	O4'-C1'-N1	5.95	112.96	108.20
53	B5	2142	A	O4'-C1'-N9	5.95	112.96	108.20
53	B5	2969	A	O4'-C1'-N9	5.95	112.96	108.20
1	AA	257	A	O4'-C1'-N9	5.95	112.96	108.20
1	AA	1263	G	O4'-C1'-N9	5.95	112.96	108.20
53	B5	364	G	O4'-C1'-N9	5.95	112.96	108.20
53	B5	1557	A	O4'-C1'-N9	5.95	112.96	108.20
1	AA	574	G	C5-C6-O6	-5.95	125.03	128.60
1	AA	578	U	O4'-C1'-N1	5.95	112.96	108.20
1	AA	876	G	N1-C6-O6	5.95	123.47	119.90
1	AA	1615	U	O4'-C1'-N1	5.95	112.96	108.20
19	A7	30	G	N9-C1'-C2'	-5.95	105.46	112.00
53	B5	552	G	N1-C6-O6	5.95	123.47	119.90
53	B5	1728	G	N1-C6-O6	5.95	123.47	119.90
53	B5	1896	A	O4'-C1'-N9	5.95	112.96	108.20
53	B5	2466	G	O4'-C1'-N9	5.95	112.96	108.20
1	AA	243	G	O4'-C1'-N9	5.94	112.95	108.20
1	AA	1570	G	C5-C6-O6	-5.94	125.03	128.60
52	B4	77	A	O4'-C1'-N9	5.94	112.95	108.20
53	B5	567	G	C5-C6-O6	-5.94	125.03	128.60
53	B5	1074	U	O4'-C1'-N1	5.94	112.95	108.20
53	B5	1326	A	O4'-C1'-N9	5.94	112.95	108.20
53	B5	1635	G	C5-C6-O6	-5.94	125.03	128.60
53	B5	1801	U	O4'-C1'-N1	5.94	112.95	108.20
1	AA	816	G	N1-C6-O6	5.94	123.46	119.90
1	AA	1123	G	C5-C6-O6	-5.94	125.04	128.60
1	AA	1496	G	C5-C6-O6	-5.94	125.03	128.60
53	B5	363	G	C5-C6-O6	-5.94	125.03	128.60
53	B5	1611	G	C5-C6-O6	-5.94	125.04	128.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	1881	A	O4'-C1'-N9	5.94	112.95	108.20
53	B5	2247	G	N1-C6-O6	5.94	123.47	119.90
1	AA	335	U	O4'-C1'-N1	5.94	112.95	108.20
52	B4	135	G	O4'-C1'-N9	5.94	112.95	108.20
53	B5	382	U	O4'-C1'-N1	5.94	112.95	108.20
53	B5	415	G	O4'-C1'-N9	5.94	112.95	108.20
53	B5	1311	G	N1-C6-O6	5.94	123.46	119.90
53	B5	2102	U	O4'-C1'-N1	5.94	112.95	108.20
53	B5	416	A	O4'-C1'-N9	5.94	112.95	108.20
52	B4	132	G	N1-C6-O6	5.93	123.46	119.90
53	B5	608	A	P-O3'-C3'	5.93	126.82	119.70
53	B5	1063	G	O4'-C1'-N9	5.93	112.95	108.20
53	B5	1763	U	O4'-C1'-N1	5.93	112.95	108.20
53	B5	2844	C	O4'-C1'-N1	5.93	112.95	108.20
1	AA	706	A	O4'-C1'-N9	5.93	112.95	108.20
1	AA	713	A	O4'-C1'-N9	5.93	112.94	108.20
1	AA	841	U	O4'-C1'-N1	5.93	112.95	108.20
1	AA	1251	G	O4'-C1'-N9	5.93	112.94	108.20
1	AA	1620	G	C5-C6-O6	-5.93	125.04	128.60
1	AA	1699	A	O4'-C1'-N9	5.93	112.95	108.20
53	B5	364	G	N1-C6-O6	5.93	123.46	119.90
53	B5	664	U	O4'-C1'-N1	5.93	112.95	108.20
1	AA	108	A	O4'-C1'-N9	5.93	112.94	108.20
1	AA	832	U	O4'-C1'-N1	5.93	112.94	108.20
1	AA	1143	G	O4'-C1'-N9	5.93	112.94	108.20
51	B3	5	G	N1-C6-O6	5.93	123.46	119.90
53	B5	875	G	O4'-C1'-N9	5.93	112.94	108.20
1	AA	1529	G	C5-C6-O6	-5.93	125.04	128.60
19	A7	70	C	C3'-C2'-C1'	-5.93	96.76	101.50
53	B5	2851	A	O4'-C1'-N9	5.93	112.94	108.20
53	B5	3174	A	O4'-C1'-N9	5.93	112.94	108.20
1	AA	284	G	N1-C6-O6	5.93	123.46	119.90
1	AA	1044	G	N1-C6-O6	5.93	123.46	119.90
53	B5	1513	G	C5-C6-O6	-5.93	125.04	128.60
53	B5	1141	C	P-O3'-C3'	5.93	126.81	119.70
1	AA	764	U	O4'-C1'-N1	5.92	112.94	108.20
1	AA	1175	G	O4'-C1'-N9	5.92	112.94	108.20
52	B4	69	U	O4'-C1'-N1	5.92	112.94	108.20
53	B5	203	G	O4'-C1'-N9	5.92	112.94	108.20
53	B5	1186	G	O4'-C1'-N9	5.92	112.94	108.20
53	B5	1313	G	C5-C6-O6	-5.92	125.05	128.60
53	B5	3049	A	C5-C6-N6	-5.92	118.96	123.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	3059	G	O4'-C1'-N9	5.92	112.94	108.20
53	B5	3220	G	C5-C6-O6	-5.92	125.05	128.60
1	AA	815	G	O4'-C1'-N9	5.92	112.94	108.20
1	AA	964	U	O4'-C1'-N1	5.92	112.94	108.20
53	B5	1029	G	O4'-C1'-N9	5.92	112.94	108.20
1	AA	365	G	C5-C6-O6	-5.92	125.05	128.60
1	AA	492	A	O4'-C1'-N9	5.92	112.94	108.20
1	AA	616	G	C5-C6-O6	-5.92	125.05	128.60
1	AA	1547	C	O3'-P-O5'	-5.92	92.75	104.00
1	AA	1716	G	O4'-C1'-N9	5.92	112.94	108.20
53	B5	1778	G	O4'-C1'-N9	5.92	112.94	108.20
53	B5	2419	A	C4-C5-C6	5.92	119.96	117.00
53	B5	24	G	O4'-C1'-N9	5.92	112.94	108.20
53	B5	718	G	N1-C6-O6	5.92	123.45	119.90
53	B5	2863	G	C1'-O4'-C4'	-5.92	105.16	109.90
53	B5	759	U	O4'-C1'-N1	5.92	112.93	108.20
53	B5	791	A	C5-C6-N6	-5.92	118.97	123.70
53	B5	854	G	N1-C6-O6	5.92	123.45	119.90
53	B5	1678	G	C5-C6-O6	-5.92	125.05	128.60
53	B5	1811	G	C5-C6-O6	-5.92	125.05	128.60
53	B5	2455	U	O4'-C1'-N1	5.92	112.94	108.20
53	B5	2775	U	O4'-C1'-N1	5.92	112.93	108.20
28	BD	50	TYR	CB-CG-CD1	-5.92	117.45	121.00
52	B4	56	G	C5-C6-O6	-5.92	125.05	128.60
53	B5	812	G	O4'-C1'-N9	5.92	112.93	108.20
53	B5	1590	G	N1-C6-O6	5.92	123.45	119.90
53	B5	3136	G	C5-C6-O6	-5.92	125.05	128.60
1	AA	1534	G	O4'-C1'-N9	5.91	112.93	108.20
52	B4	70	A	O4'-C1'-N9	5.91	112.93	108.20
53	B5	3141	A	O4'-C1'-N9	5.91	112.93	108.20
1	AA	249	U	O4'-C1'-N1	5.91	112.93	108.20
1	AA	1383	G	O4'-C1'-N9	5.91	112.93	108.20
53	B5	2937	G	O4'-C1'-N9	5.91	112.93	108.20
51	B3	65	G	C5-C6-O6	-5.91	125.05	128.60
53	B5	2609	A	O4'-C1'-N9	5.91	112.93	108.20
53	B5	219	A	C5-C6-N6	-5.91	118.97	123.70
1	AA	537	G	N1-C6-O6	5.91	123.44	119.90
1	AA	740	A	O4'-C1'-N9	5.91	112.93	108.20
1	AA	1498	C	O4'-C1'-N1	5.91	112.93	108.20
53	B5	1687	U	O4'-C1'-N1	5.91	112.93	108.20
1	AA	547	U	O4'-C1'-N1	5.91	112.92	108.20
1	AA	1294	G	O4'-C1'-N9	5.91	112.92	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1517	U	O4'-C1'-N1	5.91	112.92	108.20
1	AA	1766	G	C5-C6-O6	-5.91	125.06	128.60
53	B5	335	G	O4'-C1'-N9	5.91	112.92	108.20
53	B5	420	G	C5-C6-O6	-5.91	125.06	128.60
53	B5	916	G	C5-C6-O6	-5.91	125.06	128.60
53	B5	2587	U	O4'-C1'-N1	5.91	112.92	108.20
53	B5	2788	C	C2-N1-C1'	5.90	125.29	118.80
1	AA	1405	G	C5-C6-O6	-5.90	125.06	128.60
1	AA	1641	U	O4'-C1'-N1	5.90	112.92	108.20
53	B5	2185	G	C5-C6-O6	-5.90	125.06	128.60
1	AA	787	G	N1-C6-O6	5.90	123.44	119.90
1	AA	993	G	O4'-C1'-N9	5.90	112.92	108.20
1	AA	1349	G	O4'-C1'-N9	5.90	112.92	108.20
1	AA	1747	A	O4'-C1'-N9	5.90	112.92	108.20
53	B5	219	A	C4-C5-C6	5.90	119.95	117.00
53	B5	1357	G	O4'-C1'-N9	5.90	112.92	108.20
53	B5	2919	A	O4'-C1'-N9	5.90	112.92	108.20
1	AA	623	A	O4'-C1'-N9	5.90	112.92	108.20
1	AA	1002	G	C5-C6-O6	-5.90	125.06	128.60
1	AA	1266	G	O4'-C1'-N9	5.90	112.92	108.20
53	B5	3051	U	O4'-C1'-N1	5.90	112.92	108.20
1	AA	720	G	C5-C6-O6	-5.90	125.06	128.60
53	B5	101	G	O4'-C1'-N9	5.90	112.92	108.20
53	B5	964	G	O4'-C1'-N9	5.90	112.92	108.20
53	B5	1346	G	O4'-C1'-N9	5.90	112.92	108.20
53	B5	1434	G	N1-C6-O6	5.90	123.44	119.90
1	AA	837	G	N1-C6-O6	5.90	123.44	119.90
53	B5	3386	G	O4'-C1'-N9	5.90	112.92	108.20
1	AA	1705	A	O4'-C1'-N9	5.89	112.92	108.20
52	B4	104	A	O4'-C1'-N9	5.89	112.92	108.20
53	B5	610	G	N1-C6-O6	5.89	123.44	119.90
53	B5	1072	G	O4'-C1'-N9	5.89	112.92	108.20
53	B5	1271	A	O4'-C1'-N9	5.89	112.92	108.20
1	AA	752	A	O4'-C1'-N9	5.89	112.91	108.20
53	B5	1236	G	O4'-C1'-N9	5.89	112.91	108.20
53	B5	1473	G	O4'-C1'-N9	5.89	112.91	108.20
53	B5	2828	G	P-O3'-C3'	5.89	126.77	119.70
53	B5	2908	G	O4'-C1'-N9	5.89	112.91	108.20
1	AA	142	G	O4'-C1'-N9	5.89	112.91	108.20
1	AA	1666	G	O4'-C1'-N9	5.89	112.91	108.20
53	B5	2186	U	O4'-C1'-N1	5.89	112.91	108.20
1	AA	884	G	O4'-C1'-N9	5.89	112.91	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1694	G	O4'-C1'-N9	5.89	112.91	108.20
51	B3	63	G	C5-C6-O6	-5.89	125.07	128.60
53	B5	2131	A	C5-C6-N6	-5.89	118.99	123.70
19	A7	19	G	C6-C5-N7	-5.88	126.87	130.40
52	B4	85	G	N1-C6-O6	5.88	123.43	119.90
53	B5	837	A	O4'-C1'-N9	5.88	112.91	108.20
53	B5	1054	A	O4'-C1'-N9	5.88	112.91	108.20
53	B5	1106	G	O4'-C1'-N9	5.88	112.91	108.20
53	B5	2124	G	N1-C6-O6	5.88	123.43	119.90
53	B5	3155	U	O4'-C1'-N1	5.88	112.91	108.20
1	AA	234	G	O4'-C1'-N9	5.88	112.91	108.20
1	AA	651	G	N1-C6-O6	5.88	123.43	119.90
53	B5	1096	U	O4'-C1'-N1	5.88	112.91	108.20
1	AA	532	U	O4'-C1'-N1	5.88	112.91	108.20
1	AA	1526	U	O4'-C1'-N1	5.88	112.91	108.20
53	B5	1764	U	C2-N1-C1'	5.88	124.76	117.70
53	B5	2335	G	C5-C6-O6	-5.88	125.07	128.60
53	B5	2828	G	C5-C6-O6	-5.88	125.07	128.60
1	AA	1551	G	O4'-C1'-N9	5.88	112.90	108.20
53	B5	80	G	O4'-C1'-N9	5.88	112.90	108.20
53	B5	202	G	O4'-C1'-N9	5.88	112.90	108.20
53	B5	2280	A	O4'-C1'-N9	5.88	112.90	108.20
53	B5	3327	G	N1-C6-O6	5.88	123.43	119.90
1	AA	1346	G	N1-C6-O6	5.88	123.43	119.90
52	B4	123	G	O4'-C1'-N9	5.88	112.90	108.20
53	B5	2269	U	O4'-C1'-N1	5.88	112.90	108.20
1	AA	613	G	O4'-C1'-N9	5.88	112.90	108.20
1	AA	866	G	C3'-C2'-C1'	-5.88	96.80	101.50
1	AA	1082	G	C5-C6-O6	-5.88	125.07	128.60
1	AA	1109	G	O4'-C1'-N9	5.88	112.90	108.20
53	B5	579	G	O4'-C1'-N9	5.88	112.90	108.20
53	B5	583	G	O4'-C1'-N9	5.88	112.90	108.20
53	B5	2607	G	O4'-C1'-N9	5.88	112.90	108.20
53	B5	2895	G	O4'-C1'-N9	5.88	112.90	108.20
53	B5	219	A	O4'-C1'-N9	5.88	112.90	108.20
53	B5	1701	C	P-O3'-C3'	5.88	126.75	119.70
53	B5	2949	U	O4'-C1'-N1	5.88	112.90	108.20
1	AA	1487	U	O4'-C1'-N1	5.87	112.90	108.20
1	AA	1634	C	C4'-C3'-C2'	5.87	108.47	102.60
1	AA	1686	U	O4'-C1'-N1	5.87	112.90	108.20
53	B5	799	G	O4'-C1'-N9	5.87	112.90	108.20
53	B5	2615	G	O4'-C1'-N9	5.87	112.90	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	2784	G	C5-C6-O6	-5.87	125.08	128.60
53	B5	2786	G	C5-C6-O6	-5.87	125.08	128.60
53	B5	3247	G	O4'-C1'-N9	5.87	112.90	108.20
1	AA	751	G	O4'-C1'-N9	5.87	112.90	108.20
1	AA	1522	A	O4'-C4'-C3'	-5.87	98.13	104.00
53	B5	2140	U	O4'-C1'-N1	5.87	112.90	108.20
53	B5	2371	G	C5-C6-O6	-5.87	125.08	128.60
53	B5	2586	G	O4'-C1'-N9	5.87	112.90	108.20
53	B5	2815	G	O4'-C1'-N9	5.87	112.90	108.20
53	B5	2938	G	O4'-C1'-N9	5.87	112.90	108.20
53	B5	1371	G	C5-C6-O6	-5.87	125.08	128.60
53	B5	1455	U	O4'-C1'-N1	5.87	112.89	108.20
53	B5	3003	G	C5-C6-O6	-5.87	125.08	128.60
1	AA	284	G	O4'-C1'-N9	5.87	112.89	108.20
1	AA	1288	G	N1-C6-O6	5.87	123.42	119.90
53	B5	590	G	O4'-C1'-N9	5.87	112.89	108.20
53	B5	2558	U	O4'-C1'-N1	5.87	112.89	108.20
1	AA	956	G	N1-C6-O6	5.87	123.42	119.90
1	AA	1265	U	O4'-C1'-N1	5.87	112.89	108.20
1	AA	1293	A	O4'-C1'-N9	5.87	112.89	108.20
53	B5	424	G	C5-C6-O6	-5.87	125.08	128.60
53	B5	994	G	C5-C6-O6	-5.87	125.08	128.60
53	B5	2656	A	O4'-C1'-N9	5.87	112.89	108.20
53	B5	2978	U	O4'-C1'-N1	5.87	112.89	108.20
1	AA	516	G	N1-C6-O6	5.86	123.42	119.90
1	AA	1017	U	O4'-C1'-N1	5.86	112.89	108.20
1	AA	1720	A	O4'-C1'-N9	5.86	112.89	108.20
52	B4	49	G	C5-C6-O6	-5.86	125.08	128.60
53	B5	493	G	C5-C6-O6	-5.86	125.08	128.60
53	B5	2489	C	N3-C4-N4	5.86	122.10	118.00
1	AA	5	U	O4'-C1'-N1	5.86	112.89	108.20
1	AA	1014	U	O4'-C1'-N1	5.86	112.89	108.20
19	A7	9	A	N3-C4-C5	-5.86	122.70	126.80
53	B5	1135	A	O4'-C1'-N9	5.86	112.89	108.20
53	B5	1250	G	O4'-C1'-N9	5.86	112.89	108.20
53	B5	3045	G	O4'-C1'-N9	5.86	112.89	108.20
1	AA	165	G	O4'-C1'-N9	5.86	112.89	108.20
1	AA	498	G	O4'-C1'-N9	5.86	112.89	108.20
1	AA	771	A	O4'-C1'-N9	5.86	112.89	108.20
1	AA	853	G	O4'-C1'-N9	5.86	112.89	108.20
53	B5	566	G	C5-C6-O6	-5.86	125.08	128.60
53	B5	1374	G	O4'-C1'-N9	5.86	112.89	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	340	U	O4'-C1'-N1	5.86	112.88	108.20
1	AA	1402	G	O4'-C1'-N9	5.86	112.89	108.20
1	AA	1509	U	O4'-C1'-N1	5.86	112.88	108.20
52	B4	124	G	O4'-C1'-N9	5.86	112.89	108.20
53	B5	146	U	O4'-C1'-N1	5.86	112.89	108.20
53	B5	1852	G	C5-C6-O6	-5.86	125.09	128.60
53	B5	2728	G	O4'-C1'-N9	5.86	112.88	108.20
53	B5	3108	G	O4'-C1'-N9	5.86	112.88	108.20
1	AA	1343	A	O4'-C1'-N9	5.85	112.88	108.20
1	AA	1430	G	N1-C6-O6	5.85	123.41	119.90
53	B5	635	G	O4'-C1'-N9	5.85	112.88	108.20
53	B5	3250	U	O4'-C1'-N1	5.85	112.88	108.20
1	AA	123	G	O4'-C1'-N9	5.85	112.88	108.20
1	AA	765	G	N1-C6-O6	5.85	123.41	119.90
1	AA	824	G	O4'-C1'-N9	5.85	112.88	108.20
1	AA	1303	C	C6-N1-C2	-5.85	117.96	120.30
26	BB	156	LYS	N-CA-CB	5.85	121.13	110.60
53	B5	298	U	O4'-C1'-N1	5.85	112.88	108.20
53	B5	2199	G	C5-C6-O6	-5.85	125.09	128.60
6	AG	164	PRO	CA-N-CD	-5.85	103.31	111.50
28	BD	194	TYR	CB-CG-CD2	5.85	124.51	121.00
52	B4	9	A	O4'-C1'-N9	5.85	112.88	108.20
53	B5	1018	G	N1-C6-O6	5.85	123.41	119.90
53	B5	3031	G	C5-C6-O6	-5.85	125.09	128.60
53	B5	3291	G	N1-C6-O6	5.85	123.41	119.90
1	AA	677	G	O4'-C1'-N9	5.85	112.88	108.20
1	AA	1538	G	C5-C6-O6	-5.85	125.09	128.60
53	B5	430	U	O4'-C1'-N1	5.85	112.88	108.20
53	B5	2273	G	C5-C6-O6	-5.85	125.09	128.60
1	AA	889	U	O4'-C1'-N1	5.85	112.88	108.20
1	AA	1693	G	N1-C6-O6	5.85	123.41	119.90
53	B5	2239	G	C5-C6-O6	-5.85	125.09	128.60
53	B5	2297	U	C1'-O4'-C4'	-5.85	105.22	109.90
53	B5	2501	U	O4'-C1'-N1	5.85	112.88	108.20
1	AA	643	G	O4'-C1'-N9	5.84	112.88	108.20
19	A7	60	C	C4'-C3'-C2'	5.84	108.44	102.60
53	B5	360	G	C5-C6-O6	-5.84	125.09	128.60
53	B5	717	C	O4'-C1'-N1	5.84	112.88	108.20
53	B5	1944	U	P-O3'-C3'	5.84	126.71	119.70
1	AA	480	G	N1-C6-O6	5.84	123.41	119.90
1	AA	607	G	N1-C6-O6	5.84	123.41	119.90
1	AA	1534	G	C5-C6-O6	-5.84	125.09	128.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	1208	U	O4'-C1'-N1	5.84	112.88	108.20
1	AA	779	U	C2-N1-C1'	5.84	124.71	117.70
1	AA	836	U	O4'-C1'-N1	5.84	112.87	108.20
1	AA	840	U	O4'-C1'-N1	5.84	112.87	108.20
1	AA	1500	G	O4'-C1'-N9	5.84	112.87	108.20
53	B5	372	A	O4'-C1'-N9	5.84	112.87	108.20
53	B5	882	A	C4-C5-C6	5.84	119.92	117.00
1	AA	524	U	O4'-C1'-N1	5.84	112.87	108.20
53	B5	774	G	C5-C6-O6	-5.84	125.10	128.60
53	B5	2492	C	O4'-C1'-N1	5.84	112.87	108.20
53	B5	2878	G	O4'-C1'-N9	5.84	112.87	108.20
53	B5	3015	G	C5-C6-O6	-5.84	125.10	128.60
1	AA	287	G	O4'-C1'-N9	5.84	112.87	108.20
1	AA	1406	G	C5-C6-O6	-5.84	125.10	128.60
53	B5	1253	U	P-O3'-C3'	5.84	126.71	119.70
53	B5	2209	U	O4'-C1'-N1	5.84	112.87	108.20
1	AA	1321	G	O4'-C1'-N9	5.84	112.87	108.20
53	B5	442	G	C5-C6-O6	-5.84	125.10	128.60
53	B5	1271	A	C5-C6-N6	-5.84	119.03	123.70
19	A7	11	C	C4'-C3'-C2'	-5.83	96.77	102.60
53	B5	215	G	O4'-C1'-N9	5.83	112.87	108.20
53	B5	336	A	O4'-C1'-N9	5.83	112.87	108.20
53	B5	1169	A	O4'-C1'-N9	5.83	112.87	108.20
53	B5	2917	G	O4'-C1'-N9	5.83	112.87	108.20
1	AA	1259	G	O4'-C1'-N9	5.83	112.87	108.20
1	AA	1351	G	N1-C6-O6	5.83	123.40	119.90
1	AA	1758	G	C5-C6-O6	-5.83	125.10	128.60
53	B5	1140	G	O4'-C1'-N9	5.83	112.87	108.20
53	B5	1489	A	O4'-C1'-N9	5.83	112.87	108.20
53	B5	2953	U	O4'-C1'-N1	5.83	112.87	108.20
53	B5	2999	U	O4'-C1'-N1	5.83	112.87	108.20
53	B5	3123	A	O4'-C1'-N9	5.83	112.87	108.20
1	AA	1696	G	N1-C6-O6	5.83	123.40	119.90
19	A7	15	G	N1-C6-O6	-5.83	116.40	119.90
19	A7	62	A	C4-C5-N7	5.83	113.62	110.70
53	B5	560	G	C5-C6-O6	-5.83	125.10	128.60
53	B5	822	G	N1-C6-O6	5.83	123.40	119.90
53	B5	1465	A	O4'-C1'-N9	5.83	112.86	108.20
53	B5	2434	U	O4'-C1'-N1	5.83	112.87	108.20
53	B5	2564	G	C5-C6-O6	-5.83	125.10	128.60
53	B5	2743	A	O4'-C1'-N9	5.83	112.87	108.20
53	B5	1851	G	O4'-C1'-N9	5.83	112.86	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	3305	A	O4'-C1'-N9	5.83	112.86	108.20
52	B4	140	G	O4'-C1'-N9	5.83	112.86	108.20
53	B5	353	G	P-O3'-C3'	5.83	126.69	119.70
53	B5	1091	A	O4'-C1'-N9	5.83	112.86	108.20
53	B5	2435	G	C5-C6-O6	-5.83	125.10	128.60
53	B5	2914	G	O4'-C1'-N9	5.83	112.86	108.20
53	B5	3325	G	N1-C6-O6	5.83	123.40	119.90
53	B5	3345	G	O4'-C1'-N9	5.83	112.86	108.20
53	B5	53	G	N1-C6-O6	5.83	123.39	119.90
53	B5	827	A	C5-C6-N6	-5.83	119.04	123.70
53	B5	828	A	C5-C6-N6	-5.83	119.04	123.70
53	B5	926	A	C5-C6-N6	-5.83	119.04	123.70
53	B5	1171	G	C5-C6-O6	-5.83	125.11	128.60
53	B5	1247	U	O4'-C1'-N1	5.83	112.86	108.20
53	B5	1438	U	O4'-C1'-N1	5.83	112.86	108.20
53	B5	1469	C	C2-N1-C1'	5.83	125.21	118.80
53	B5	2262	A	O4'-C1'-N9	5.83	112.86	108.20
53	B5	306	A	O4'-C1'-N9	5.82	112.86	108.20
53	B5	538	G	C5-C6-O6	-5.82	125.11	128.60
53	B5	1052	U	O4'-C1'-N1	5.82	112.86	108.20
53	B5	1480	G	O4'-C1'-N9	5.82	112.86	108.20
1	AA	733	A	O4'-C1'-N9	5.82	112.86	108.20
19	A7	9	A	C5-N7-C8	5.82	106.81	103.90
51	B3	86	G	O4'-C1'-N9	5.82	112.86	108.20
53	B5	2559	A	O4'-C1'-N9	5.82	112.86	108.20
1	AA	1780	A	O4'-C1'-N9	5.82	112.86	108.20
53	B5	235	A	O4'-C1'-N9	5.82	112.86	108.20
53	B5	3015	G	O4'-C1'-N9	5.82	112.86	108.20
1	AA	1612	A	P-O5'-C5'	-5.82	111.59	120.90
1	AA	285	G	O4'-C1'-N9	5.82	112.86	108.20
1	AA	1481	A	O4'-C1'-N9	5.82	112.85	108.20
53	B5	1011	A	O4'-C1'-N9	5.82	112.85	108.20
53	B5	1321	G	O4'-C1'-N9	5.82	112.85	108.20
53	B5	2315	G	N1-C6-O6	5.82	123.39	119.90
53	B5	2462	A	O4'-C1'-N9	5.82	112.85	108.20
53	B5	2639	G	C5-C6-O6	-5.82	125.11	128.60
53	B5	3345	G	C5-C6-O6	-5.82	125.11	128.60
1	AA	400	A	O4'-C1'-N9	5.82	112.85	108.20
1	AA	751	G	N1-C6-O6	5.82	123.39	119.90
1	AA	846	G	C5-C6-O6	-5.82	125.11	128.60
1	AA	1278	U	O4'-C1'-N1	5.82	112.85	108.20
53	B5	1323	G	N1-C6-O6	5.82	123.39	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	631	G	C5-C6-O6	-5.81	125.11	128.60
1	AA	1066	A	O4'-C1'-N9	5.81	112.85	108.20
53	B5	2396	G	O4'-C1'-N9	5.81	112.85	108.20
53	B5	2457	G	O4'-C1'-N9	5.81	112.85	108.20
1	AA	507	U	O4'-C1'-N1	5.81	112.85	108.20
1	AA	1086	U	O4'-C1'-N1	5.81	112.85	108.20
53	B5	1270	A	O4'-C1'-N9	5.81	112.85	108.20
53	B5	1362	G	C5-C6-O6	-5.81	125.11	128.60
53	B5	845	G	P-O3'-C3'	5.81	126.67	119.70
1	AA	220	A	O4'-C1'-N9	5.81	112.85	108.20
1	AA	734	A	O4'-C1'-N9	5.81	112.85	108.20
1	AA	1478	G	O4'-C1'-N9	5.81	112.85	108.20
1	AA	1657	A	O4'-C1'-N9	5.81	112.85	108.20
19	A7	45	G	C8-N9-C4	-5.81	104.08	106.40
53	B5	1226	G	O4'-C1'-N9	5.81	112.85	108.20
53	B5	2311	G	N1-C6-O6	5.81	123.39	119.90
1	AA	875	G	C5-C6-O6	-5.81	125.12	128.60
53	B5	30	G	C5-C6-O6	-5.81	125.11	128.60
1	AA	1664	U	O4'-C1'-N1	5.81	112.84	108.20
53	B5	335	G	C5-C6-O6	-5.81	125.12	128.60
1	AA	206	A	O4'-C1'-N9	5.80	112.84	108.20
1	AA	606	A	O4'-C1'-N9	5.80	112.84	108.20
1	AA	689	G	N1-C6-O6	5.80	123.38	119.90
53	B5	213	G	C5-C6-O6	-5.80	125.12	128.60
53	B5	2661	G	N1-C6-O6	5.80	123.38	119.90
53	B5	3013	U	O4'-C1'-N1	5.80	112.84	108.20
53	B5	3363	U	P-O3'-C3'	5.80	126.66	119.70
53	B5	475	G	C5-C6-O6	-5.80	125.12	128.60
53	B5	2917	G	N1-C6-O6	5.80	123.38	119.90
1	AA	535	A	O4'-C1'-N9	5.80	112.84	108.20
17	AR	164	ASP	N-CA-C	-5.80	95.33	111.00
53	B5	269	G	C5-C6-O6	-5.80	125.12	128.60
53	B5	1375	G	C5-C6-O6	-5.80	125.12	128.60
53	B5	1742	U	O4'-C1'-N1	5.80	112.84	108.20
1	AA	3	U	O4'-C1'-N1	5.80	112.84	108.20
1	AA	236	A	O4'-C1'-N9	5.80	112.84	108.20
53	B5	1753	G	C5-C6-O6	-5.80	125.12	128.60
53	B5	2161	G	N1-C6-O6	5.80	123.38	119.90
19	A7	51	G	N1-C2-N3	5.80	127.38	123.90
53	B5	337	G	O4'-C1'-N9	5.80	112.84	108.20
53	B5	737	G	O4'-C1'-N9	5.80	112.84	108.20
53	B5	2748	A	C5-C6-N6	-5.80	119.06	123.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	3320	A	C5-C6-N6	-5.80	119.06	123.70
53	B5	2766	U	O4'-C1'-N1	5.79	112.84	108.20
53	B5	2995	A	O4'-C1'-N9	5.79	112.84	108.20
1	AA	396	G	C5-C6-O6	-5.79	125.12	128.60
1	AA	1364	G	O4'-C1'-N9	5.79	112.83	108.20
1	AA	1430	G	P-O3'-C3'	5.79	126.65	119.70
1	AA	1678	G	N1-C6-O6	5.79	123.38	119.90
53	B5	1790	G	O4'-C1'-N9	5.79	112.83	108.20
53	B5	2968	G	O4'-C1'-N9	5.79	112.83	108.20
1	AA	1098	G	C5-C6-O6	-5.79	125.12	128.60
19	A7	43	G	C4-C5-N7	-5.79	108.48	110.80
19	A7	60	C	N1-C2-N3	5.79	123.25	119.20
51	B3	106	G	C5-C6-O6	-5.79	125.12	128.60
53	B5	1474	A	O4'-C1'-N9	5.79	112.83	108.20
53	B5	1581	C	N3-C4-N4	5.79	122.06	118.00
53	B5	2972	G	C5-C6-O6	-5.79	125.12	128.60
53	B5	3198	U	O4'-C1'-N1	5.79	112.83	108.20
1	AA	1588	G	C5-C6-O6	-5.79	125.13	128.60
1	AA	1535	C	O4'-C1'-N1	5.79	112.83	108.20
53	B5	766	U	O4'-C1'-N1	5.79	112.83	108.20
1	AA	533	U	O4'-C1'-N1	5.79	112.83	108.20
1	AA	1045	G	N1-C6-O6	5.79	123.37	119.90
1	AA	1586	G	C5-C6-O6	-5.79	125.13	128.60
19	A7	14	A	C5-C6-N1	5.79	120.59	117.70
1	AA	37	U	O4'-C1'-N1	5.79	112.83	108.20
1	AA	877	G	N1-C6-O6	5.79	123.37	119.90
53	B5	1635	G	O4'-C1'-N9	5.79	112.83	108.20
53	B5	1778	G	C5-C6-O6	-5.79	125.13	128.60
1	AA	308	C	O4'-C1'-N1	5.78	112.83	108.20
1	AA	1582	G	C5-C6-O6	-5.78	125.13	128.60
1	AA	1670	G	N1-C6-O6	5.78	123.37	119.90
53	B5	534	U	O4'-C1'-N1	5.78	112.83	108.20
53	B5	1709	C	O4'-C1'-N1	5.78	112.83	108.20
53	B5	2240	G	C5-C6-O6	-5.78	125.13	128.60
53	B5	1443	G	C5-C6-O6	-5.78	125.13	128.60
1	AA	311	U	O4'-C1'-N1	5.78	112.82	108.20
1	AA	602	U	O4'-C1'-N1	5.78	112.82	108.20
1	AA	1034	G	O4'-C1'-N9	5.78	112.82	108.20
53	B5	1048	A	C5-C6-N6	-5.78	119.08	123.70
53	B5	1591	G	O4'-C1'-N9	5.78	112.83	108.20
53	B5	2912	G	O4'-C1'-N9	5.78	112.83	108.20
53	B5	3223	A	O4'-C1'-N9	5.78	112.83	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	3241	G	O4'-C1'-N9	5.78	112.82	108.20
19	A7	70	C	N1-C1'-C2'	-5.78	105.64	112.00
53	B5	24	G	C5-C6-O6	-5.78	125.13	128.60
1	AA	969	A	O4'-C1'-N9	5.78	112.82	108.20
53	B5	273	A	O4'-C1'-N9	5.78	112.82	108.20
53	B5	2952	G	N1-C6-O6	5.78	123.37	119.90
1	AA	228	G	O4'-C1'-N9	5.78	112.82	108.20
1	AA	1489	U	O4'-C1'-N1	5.78	112.82	108.20
1	AA	1593	U	P-O3'-C3'	5.78	126.63	119.70
53	B5	2902	A	O4'-C1'-N9	5.78	112.82	108.20
53	B5	3229	G	C5-C6-O6	-5.78	125.13	128.60
1	AA	979	G	O4'-C1'-N9	5.77	112.82	108.20
52	B4	149	A	C5-C6-N6	-5.77	119.08	123.70
53	B5	842	G	C5-C6-O6	-5.77	125.14	128.60
53	B5	1537	A	O4'-C1'-N9	5.77	112.82	108.20
1	AA	428	A	O4'-C1'-N9	5.77	112.82	108.20
1	AA	984	G	O4'-C1'-N9	5.77	112.82	108.20
19	A7	42	G	O4'-C1'-N9	5.77	112.82	108.20
53	B5	2171	G	O4'-C1'-N9	5.77	112.82	108.20
53	B5	3087	A	C4-C5-C6	5.77	119.89	117.00
53	B5	3216	U	O4'-C1'-N1	5.77	112.82	108.20
1	AA	373	G	C5-C6-O6	-5.77	125.14	128.60
19	A7	73	A	N1-C6-N6	-5.77	115.14	118.60
53	B5	53	G	O4'-C1'-N9	5.77	112.82	108.20
1	AA	207	U	O4'-C1'-N1	5.77	112.82	108.20
1	AA	1050	G	N1-C6-O6	5.77	123.36	119.90
12	AM	85	PHE	CB-CG-CD1	5.77	124.84	120.80
19	A7	53	G	C8-N9-C4	-5.77	104.09	106.40
1	AA	1237	G	O4'-C1'-N9	5.77	112.81	108.20
1	AA	1679	A	O4'-C1'-N9	5.77	112.81	108.20
1	AA	1719	A	O4'-C1'-N9	5.77	112.81	108.20
52	B4	144	G	O4'-C1'-N9	5.77	112.81	108.20
53	B5	170	G	N1-C6-O6	5.77	123.36	119.90
53	B5	909	G	O4'-C1'-N9	5.77	112.81	108.20
53	B5	2194	G	C5-C6-O6	-5.77	125.14	128.60
18	AT	33	TYR	CB-CG-CD2	-5.77	117.54	121.00
53	B5	2278	C	C2-N1-C1'	5.77	125.14	118.80
1	AA	432	G	C5-C6-O6	-5.76	125.14	128.60
51	B3	30	G	O4'-C1'-N9	5.76	112.81	108.20
53	B5	710	A	C5-C6-N6	-5.76	119.09	123.70
1	AA	1678	G	O4'-C1'-N9	5.76	112.81	108.20
19	A7	4	G	C1'-O4'-C4'	5.76	114.51	109.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	34	G	N1-C6-O6	5.76	123.36	119.90
1	AA	558	U	O4'-C1'-N1	5.76	112.81	108.20
1	AA	1291	G	O4'-C1'-N9	5.76	112.81	108.20
1	AA	1733	U	O4'-C1'-N1	5.76	112.81	108.20
51	B3	96	G	C5-C6-O6	-5.76	125.14	128.60
53	B5	307	A	O4'-C1'-N9	5.76	112.81	108.20
53	B5	1915	A	C4-C5-C6	5.76	119.88	117.00
53	B5	2563	G	C5-C6-O6	-5.76	125.14	128.60
1	AA	1382	G	N1-C6-O6	5.76	123.36	119.90
19	A7	44	A	C5-C6-N6	5.76	128.31	123.70
53	B5	571	U	O4'-C1'-N1	5.76	112.81	108.20
53	B5	1206	G	C5-C6-O6	-5.76	125.14	128.60
53	B5	2622	C	N3-C4-N4	5.76	122.03	118.00
53	B5	2993	G	C5-C6-O6	-5.76	125.14	128.60
1	AA	1751	A	O4'-C1'-N9	5.76	112.81	108.20
53	B5	195	U	O4'-C1'-N1	5.76	112.81	108.20
53	B5	702	C	N3-C4-N4	5.76	122.03	118.00
53	B5	1799	A	O4'-C1'-N9	5.76	112.81	108.20
1	AA	143	G	O4'-C1'-N9	5.76	112.81	108.20
1	AA	1560	G	O4'-C1'-N9	5.76	112.81	108.20
53	B5	1010	G	C5-C6-O6	-5.76	125.15	128.60
53	B5	1310	G	C5-C6-O6	-5.76	125.15	128.60
53	B5	1604	G	O4'-C1'-N9	5.76	112.81	108.20
53	B5	2606	G	O4'-C1'-N9	5.76	112.81	108.20
53	B5	3270	A	O4'-C1'-N9	5.76	112.81	108.20
1	AA	359	A	O4'-C1'-N9	5.75	112.80	108.20
1	AA	1556	U	O4'-C1'-N1	5.75	112.80	108.20
53	B5	1882	G	C5-C6-O6	-5.75	125.15	128.60
53	B5	2169	G	O4'-C1'-N9	5.75	112.80	108.20
1	AA	11	A	O4'-C1'-N9	5.75	112.80	108.20
1	AA	539	G	O4'-C1'-N9	5.75	112.80	108.20
1	AA	690	G	O4'-C1'-N9	5.75	112.80	108.20
1	AA	1671	G	O4'-C1'-N9	5.75	112.80	108.20
53	B5	755	A	O4'-C1'-N9	5.75	112.80	108.20
53	B5	1039	U	O4'-C1'-N1	5.75	112.80	108.20
53	B5	1743	G	O4'-C1'-N9	5.75	112.80	108.20
51	B3	59	G	O4'-C1'-N9	5.75	112.80	108.20
53	B5	1010	G	O4'-C1'-N9	5.75	112.80	108.20
1	AA	1229	G	O4'-C1'-N9	5.75	112.80	108.20
53	B5	714	G	O4'-C1'-N9	5.75	112.80	108.20
53	B5	2679	A	C5-C6-N6	-5.75	119.10	123.70
53	B5	2830	G	C5-C6-O6	-5.75	125.15	128.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1122	A	C5-C6-N6	-5.75	119.10	123.70
1	AA	1173	G	C5-C6-O6	-5.75	125.15	128.60
1	AA	1400	C	N3-C4-N4	5.75	122.02	118.00
1	AA	1574	A	O4'-C1'-N9	5.75	112.80	108.20
53	B5	1751	G	C5-C6-O6	-5.75	125.15	128.60
53	B5	2721	A	O4'-C1'-N9	5.75	112.80	108.20
1	AA	1403	A	O4'-C1'-N9	5.75	112.80	108.20
1	AA	1767	U	O4'-C1'-N1	5.75	112.80	108.20
53	B5	1142	G	C5-C6-O6	-5.75	125.15	128.60
53	B5	1734	G	C5-C6-O6	-5.75	125.15	128.60
53	B5	2614	G	O4'-C1'-N9	5.75	112.80	108.20
53	B5	366	A	O4'-C1'-N9	5.75	112.80	108.20
53	B5	1766	G	O4'-C1'-N9	5.75	112.80	108.20
1	AA	8	U	O4'-C1'-N1	5.74	112.80	108.20
1	AA	175	G	C5-C6-O6	-5.74	125.15	128.60
1	AA	362	G	C5-C6-O6	-5.74	125.15	128.60
1	AA	855	A	C5-C6-N6	-5.74	119.11	123.70
1	AA	1423	C	O4'-C1'-N1	5.74	112.80	108.20
53	B5	1638	A	O4'-C1'-N9	5.74	112.80	108.20
53	B5	1509	A	O4'-C1'-N9	5.74	112.79	108.20
1	AA	153	G	C5-C6-O6	-5.74	125.16	128.60
1	AA	1665	A	O4'-C1'-N9	5.74	112.79	108.20
53	B5	1517	G	O4'-C1'-N9	5.74	112.79	108.20
53	B5	3170	A	O4'-C1'-N9	5.74	112.79	108.20
53	B5	3373	U	O4'-C1'-N1	5.74	112.79	108.20
1	AA	996	G	O4'-C1'-N9	5.74	112.79	108.20
1	AA	1238	A	O4'-C1'-N9	5.74	112.79	108.20
53	B5	182	U	O4'-C1'-N1	5.74	112.79	108.20
53	B5	662	U	O4'-C1'-N1	5.74	112.79	108.20
53	B5	1220	U	O4'-C1'-N1	5.74	112.79	108.20
53	B5	450	G	C5-C6-O6	-5.74	125.16	128.60
53	B5	505	G	C5-C6-O6	-5.74	125.16	128.60
53	B5	1529	A	O4'-C1'-N9	5.74	112.79	108.20
53	B5	2802	A	O4'-C1'-N9	5.74	112.79	108.20
53	B5	2864	A	O4'-C1'-N9	5.74	112.79	108.20
1	AA	168	A	O4'-C1'-N9	5.74	112.79	108.20
53	B5	75	G	C5-C6-O6	-5.74	125.16	128.60
53	B5	227	G	C5-C6-O6	-5.74	125.16	128.60
53	B5	1087	G	O4'-C1'-N9	5.74	112.79	108.20
53	B5	2160	G	N1-C6-O6	5.74	123.34	119.90
53	B5	2295	A	P-O5'-C5'	-5.74	111.72	120.90
53	B5	2347	U	O4'-C1'-N1	5.74	112.79	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1391	G	O4'-C1'-N9	5.73	112.79	108.20
53	B5	443	G	C5-C6-O6	-5.73	125.16	128.60
53	B5	3290	G	O4'-C1'-N9	5.73	112.79	108.20
1	AA	761	G	O4'-C1'-N9	5.73	112.79	108.20
53	B5	566	G	O4'-C1'-N9	5.73	112.79	108.20
53	B5	1259	A	O4'-C1'-N9	5.73	112.79	108.20
53	B5	1381	A	O4'-C1'-N9	5.73	112.79	108.20
53	B5	2738	A	O4'-C1'-N9	5.73	112.79	108.20
1	AA	1052	U	O4'-C1'-N1	5.73	112.78	108.20
1	AA	1591	A	O4'-C1'-N9	5.73	112.78	108.20
53	B5	189	G	N1-C6-O6	5.73	123.34	119.90
53	B5	229	G	P-O3'-C3'	5.73	126.58	119.70
53	B5	993	A	O4'-C1'-N9	5.73	112.78	108.20
53	B5	1145	G	O4'-C1'-N9	5.73	112.78	108.20
53	B5	1658	G	O4'-C1'-N9	5.73	112.78	108.20
53	B5	1724	U	O4'-C1'-N1	5.73	112.78	108.20
53	B5	3147	G	O4'-C1'-N9	5.73	112.78	108.20
1	AA	281	G	C5-C6-O6	-5.73	125.16	128.60
53	B5	220	G	C5-C6-O6	-5.73	125.16	128.60
53	B5	1676	A	O4'-C1'-N9	5.73	112.78	108.20
53	B5	1770	G	O4'-C1'-N9	5.73	112.78	108.20
53	B5	2699	G	O4'-C1'-N9	5.73	112.78	108.20
1	AA	324	U	O4'-C1'-N1	5.73	112.78	108.20
1	AA	1413	G	O4'-C1'-N9	5.73	112.78	108.20
53	B5	705	A	O4'-C1'-N9	5.73	112.78	108.20
53	B5	779	G	O4'-C1'-N9	5.73	112.78	108.20
1	AA	722	G	O4'-C1'-N9	5.73	112.78	108.20
53	B5	26	A	C5-C6-N6	-5.73	119.12	123.70
53	B5	513	G	O4'-C1'-N9	5.73	112.78	108.20
53	B5	1699	A	O4'-C1'-N9	5.73	112.78	108.20
53	B5	1817	G	O4'-C1'-N9	5.73	112.78	108.20
1	AA	222	A	O4'-C1'-N9	5.72	112.78	108.20
1	AA	326	G	C5-C6-O6	-5.72	125.17	128.60
1	AA	1460	G	C5-C6-O6	-5.72	125.17	128.60
19	A7	6	U	N1-C2-N3	5.72	118.33	114.90
53	B5	846	A	O4'-C1'-N9	5.72	112.78	108.20
53	B5	1590	G	O4'-C1'-N9	5.72	112.78	108.20
53	B5	2271	A	O4'-C1'-N9	5.72	112.78	108.20
1	AA	662	U	O4'-C1'-N1	5.72	112.78	108.20
53	B5	176	G	C5-C6-O6	-5.72	125.17	128.60
53	B5	2865	U	O4'-C1'-N1	5.72	112.78	108.20
1	AA	182	A	O4'-C1'-N9	5.72	112.78	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	228	G	C5-C6-O6	-5.72	125.17	128.60
1	AA	732	G	N1-C6-O6	5.72	123.33	119.90
1	AA	1522	A	O5'-P-OP1	-5.72	100.55	105.70
19	A7	29	A	C5-C6-N1	5.72	120.56	117.70
53	B5	2273	G	O4'-C1'-N9	5.72	112.78	108.20
53	B5	2457	G	C5-C6-O6	-5.72	125.17	128.60
53	B5	2778	G	C5-C6-O6	-5.72	125.17	128.60
53	B5	1444	G	C5-C6-O6	-5.72	125.17	128.60
1	AA	1178	U	O4'-C1'-N1	5.72	112.77	108.20
1	AA	1323	A	O4'-C1'-N9	5.72	112.78	108.20
1	AA	1712	A	O4'-C1'-N9	5.72	112.77	108.20
53	B5	701	G	C5-C6-O6	-5.72	125.17	128.60
1	AA	464	A	O4'-C1'-N9	5.72	112.77	108.20
1	AA	552	G	C5-C6-O6	-5.72	125.17	128.60
1	AA	937	G	C5-C6-O6	-5.72	125.17	128.60
1	AA	1778	G	C5-C6-O6	-5.72	125.17	128.60
52	B4	41	A	C5-C6-N6	-5.72	119.13	123.70
53	B5	213	G	O4'-C1'-N9	5.72	112.77	108.20
53	B5	483	G	O4'-C1'-N9	5.72	112.77	108.20
53	B5	722	G	O4'-C1'-N9	5.72	112.77	108.20
53	B5	2688	U	O4'-C1'-N1	5.72	112.77	108.20
53	B5	3033	A	O4'-C1'-N9	5.72	112.77	108.20
1	AA	95	G	O4'-C1'-N9	5.71	112.77	108.20
1	AA	386	G	C5-C6-O6	-5.71	125.17	128.60
1	AA	1388	A	O4'-C1'-N9	5.71	112.77	108.20
19	A7	21	A	N1-C6-N6	-5.71	115.17	118.60
53	B5	234	G	O4'-C1'-N9	5.71	112.77	108.20
1	AA	520	A	O4'-C1'-N9	5.71	112.77	108.20
53	B5	1726	C	O4'-C1'-N1	5.71	112.77	108.20
53	B5	2529	A	O4'-C1'-N9	5.71	112.77	108.20
1	AA	293	U	O4'-C1'-N1	5.71	112.77	108.20
1	AA	471	A	O4'-C1'-N9	5.71	112.77	108.20
1	AA	498	G	N1-C6-O6	5.71	123.33	119.90
1	AA	775	G	N1-C6-O6	5.71	123.33	119.90
19	A7	5	A	C5'-C4'-O4'	5.71	115.95	109.10
53	B5	426	G	O4'-C1'-N9	5.71	112.77	108.20
53	B5	1149	G	C5-C6-O6	-5.71	125.17	128.60
53	B5	2381	G	C5-C6-O6	-5.71	125.17	128.60
19	A7	30	G	C4-C5-N7	-5.71	108.52	110.80
53	B5	466	G	N1-C6-O6	5.71	123.33	119.90
53	B5	926	A	C4-C5-C6	5.71	119.86	117.00
53	B5	3049	A	O4'-C1'-N9	5.71	112.77	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	458	G	O4'-C1'-N9	5.71	112.77	108.20
1	AA	630	A	O4'-C1'-N9	5.71	112.77	108.20
1	AA	1345	A	O4'-C1'-N9	5.71	112.77	108.20
50	BZ	46	LYS	N-CA-CB	5.71	120.88	110.60
53	B5	96	G	O4'-C1'-N9	5.71	112.77	108.20
53	B5	494	G	C5-C6-O6	-5.71	125.17	128.60
53	B5	1258	U	O4'-C1'-N1	5.71	112.77	108.20
53	B5	2164	A	O4'-C1'-N9	5.71	112.77	108.20
53	B5	2414	G	C5-C6-O6	-5.71	125.17	128.60
53	B5	3157	U	O4'-C1'-N1	5.71	112.77	108.20
1	AA	721	U	O4'-C1'-N1	5.71	112.77	108.20
1	AA	1236	U	O4'-C1'-N1	5.71	112.77	108.20
53	B5	551	A	O4'-C1'-N9	5.71	112.77	108.20
53	B5	1318	A	O4'-C1'-N9	5.71	112.77	108.20
53	B5	1441	G	P-O3'-C3'	5.71	126.55	119.70
53	B5	1823	A	O4'-C1'-N9	5.71	112.76	108.20
53	B5	1935	G	N1-C6-O6	5.71	123.32	119.90
53	B5	2658	G	C5-C6-O6	-5.71	125.18	128.60
53	B5	2662	G	N1-C6-O6	5.71	123.32	119.90
1	AA	290	G	O4'-C1'-N9	5.70	112.76	108.20
1	AA	389	G	N1-C6-O6	5.70	123.32	119.90
1	AA	1758	G	O4'-C1'-N9	5.70	112.76	108.20
53	B5	318	A	C5-C6-N6	-5.70	119.14	123.70
53	B5	1526	U	O4'-C1'-N1	5.70	112.76	108.20
1	AA	515	A	O4'-C1'-N9	5.70	112.76	108.20
1	AA	1317	U	O4'-C1'-N1	5.70	112.76	108.20
51	B3	45	A	O4'-C1'-N9	5.70	112.76	108.20
53	B5	162	G	N1-C6-O6	5.70	123.32	119.90
53	B5	478	A	O4'-C1'-N9	5.70	112.76	108.20
53	B5	608	A	C5-C6-N6	-5.70	119.14	123.70
53	B5	2295	A	C5-C6-N6	-5.70	119.14	123.70
1	AA	125	U	O4'-C1'-N1	5.70	112.76	108.20
1	AA	577	G	C5-C6-O6	-5.70	125.18	128.60
1	AA	931	U	O4'-C1'-N1	5.70	112.76	108.20
1	AA	1149	A	O4'-C1'-N9	5.70	112.76	108.20
1	AA	1338	A	C5-C6-N6	-5.70	119.14	123.70
53	B5	241	G	N1-C6-O6	5.70	123.32	119.90
53	B5	800	G	C4-N9-C1'	5.70	133.91	126.50
53	B5	1333	C	N3-C4-C5	-5.70	119.62	121.90
53	B5	1493	G	O4'-C1'-N9	5.70	112.76	108.20
53	B5	1728	G	O4'-C1'-N9	5.70	112.76	108.20
53	B5	2279	A	C5-C6-N6	-5.70	119.14	123.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	3213	G	O4'-C1'-N9	5.70	112.76	108.20
1	AA	763	G	C5-C6-O6	-5.70	125.18	128.60
1	AA	1524	A	O4'-C1'-N9	5.70	112.76	108.20
52	B4	79	A	O4'-C1'-N9	5.70	112.76	108.20
53	B5	3042	U	O4'-C1'-N1	5.70	112.76	108.20
1	AA	100	A	C5-C6-N6	-5.70	119.14	123.70
1	AA	488	G	O4'-C1'-N9	5.70	112.76	108.20
19	A7	70	C	N1-C2-O2	5.70	122.32	118.90
52	B4	60	U	C5'-C4'-C3'	5.70	125.11	116.00
53	B5	1019	G	O4'-C1'-N9	5.70	112.76	108.20
53	B5	1626	U	O4'-C1'-N1	5.70	112.76	108.20
53	B5	1788	C	N3-C4-N4	5.70	121.99	118.00
1	AA	255	U	O4'-C1'-N1	5.69	112.75	108.20
1	AA	804	A	O4'-C1'-N9	5.69	112.75	108.20
53	B5	1083	G	O4'-C1'-N9	5.69	112.76	108.20
53	B5	1859	A	O4'-C1'-N9	5.69	112.75	108.20
53	B5	1734	G	O4'-C1'-N9	5.69	112.75	108.20
53	B5	2316	G	N1-C6-O6	5.69	123.31	119.90
53	B5	2468	A	P-O3'-C3'	5.69	126.53	119.70
53	B5	2877	G	O4'-C1'-N9	5.69	112.75	108.20
1	AA	30	G	O4'-C1'-N9	5.69	112.75	108.20
1	AA	1585	A	C4-C5-C6	5.69	119.84	117.00
53	B5	209	A	O4'-C1'-N9	5.69	112.75	108.20
53	B5	1268	G	O4'-C1'-N9	5.69	112.75	108.20
53	B5	1667	A	P-O3'-C3'	5.69	126.53	119.70
1	AA	892	A	O4'-C1'-N9	5.69	112.75	108.20
53	B5	476	G	C5-C6-O6	-5.69	125.19	128.60
53	B5	554	A	O4'-C1'-N9	5.69	112.75	108.20
53	B5	999	G	C5-C6-O6	-5.69	125.19	128.60
53	B5	1485	G	O4'-C1'-N9	5.69	112.75	108.20
53	B5	1921	A	O4'-C1'-N9	5.69	112.75	108.20
53	B5	2801	A	O4'-C1'-N9	5.69	112.75	108.20
1	AA	16	G	C5-C6-O6	-5.69	125.19	128.60
1	AA	357	G	N1-C6-O6	5.69	123.31	119.90
1	AA	1305	G	O4'-C1'-N9	5.69	112.75	108.20
52	B4	63	G	C5-C6-O6	-5.69	125.19	128.60
53	B5	1244	A	C1'-O4'-C4'	-5.69	105.35	109.90
53	B5	1332	A	C4-C5-C6	5.69	119.84	117.00
53	B5	763	G	O4'-C1'-N9	5.68	112.75	108.20
53	B5	1766	G	C5-C6-O6	-5.68	125.19	128.60
53	B5	2895	G	C5-C6-O6	-5.68	125.19	128.60
1	AA	391	A	O4'-C1'-N9	5.68	112.75	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	720	G	O4'-C1'-N9	5.68	112.75	108.20
1	AA	788	A	O4'-C1'-N9	5.68	112.75	108.20
53	B5	547	G	C5-C6-O6	-5.68	125.19	128.60
53	B5	876	A	O4'-C1'-N9	5.68	112.75	108.20
53	B5	1669	C	N3-C4-N4	5.68	121.98	118.00
53	B5	2519	A	O4'-C1'-N9	5.68	112.75	108.20
1	AA	789	A	O4'-C1'-N9	5.68	112.75	108.20
53	B5	798	G	O4'-C1'-N9	5.68	112.75	108.20
53	B5	2375	G	C5-C6-O6	-5.68	125.19	128.60
1	AA	362	G	O4'-C1'-N9	5.68	112.74	108.20
1	AA	508	U	O4'-C1'-N1	5.68	112.74	108.20
1	AA	634	G	O4'-C1'-N9	5.68	112.74	108.20
1	AA	1620	G	O4'-C1'-N9	5.68	112.74	108.20
29	BE	142	PHE	CB-CG-CD1	5.68	124.78	120.80
53	B5	8	C	O4'-C1'-N1	5.68	112.74	108.20
53	B5	241	G	O4'-C1'-N9	5.68	112.74	108.20
53	B5	3214	U	O4'-C1'-N1	5.68	112.74	108.20
53	B5	3286	G	O4'-C1'-N9	5.68	112.74	108.20
1	AA	867	G	C2'-C3'-O3'	-5.68	97.01	109.50
1	AA	1734	G	N1-C6-O6	5.68	123.31	119.90
53	B5	684	G	O4'-C1'-N9	5.68	112.74	108.20
1	AA	486	G	N1-C6-O6	5.68	123.31	119.90
1	AA	951	A	O4'-C1'-N9	5.68	112.74	108.20
19	A7	64	A	C5-N7-C8	5.68	106.74	103.90
53	B5	542	G	O4'-C1'-N9	5.68	112.74	108.20
53	B5	617	G	N1-C6-O6	5.68	123.31	119.90
53	B5	1414	G	O4'-C1'-N9	5.68	112.74	108.20
53	B5	2440	G	O4'-C1'-N9	5.68	112.74	108.20
1	AA	291	G	O4'-C1'-N9	5.67	112.74	108.20
1	AA	838	G	C5-C6-O6	-5.67	125.19	128.60
53	B5	389	A	O4'-C1'-N9	5.67	112.74	108.20
53	B5	740	G	C5-C6-O6	-5.67	125.19	128.60
53	B5	1112	A	C5-C6-N6	-5.67	119.16	123.70
53	B5	2463	G	N1-C6-O6	5.67	123.30	119.90
53	B5	2491	A	O4'-C1'-N9	5.67	112.74	108.20
53	B5	3331	U	O4'-C1'-N1	5.67	112.74	108.20
1	AA	60	U	O4'-C1'-N1	5.67	112.74	108.20
53	B5	1813	A	C5-C6-N6	-5.67	119.16	123.70
1	AA	756	A	O4'-C1'-N9	5.67	112.74	108.20
1	AA	1238	A	C4-C5-C6	5.67	119.84	117.00
53	B5	1059	G	C5-C6-O6	-5.67	125.20	128.60
53	B5	1373	A	O4'-C1'-N9	5.67	112.74	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	2897	A	C5-C6-N6	-5.67	119.16	123.70
50	BZ	43	TYR	CB-CG-CD1	5.67	124.40	121.00
53	B5	70	A	O4'-C1'-N9	5.67	112.74	108.20
53	B5	2734	A	O4'-C1'-N9	5.67	112.74	108.20
53	B5	2933	A	O4'-C1'-N9	5.67	112.74	108.20
1	AA	66	U	O4'-C1'-N1	5.67	112.74	108.20
1	AA	1635	C	C4'-C3'-C2'	-5.67	96.93	102.60
19	A7	3	G	N9-C4-C5	5.67	107.67	105.40
53	B5	1889	G	O4'-C1'-N9	5.67	112.73	108.20
53	B5	2749	G	C5-C6-O6	-5.67	125.20	128.60
52	B4	63	G	O4'-C1'-N9	5.67	112.73	108.20
53	B5	736	A	O4'-C1'-N9	5.67	112.73	108.20
53	B5	1072	G	C5-C6-O6	-5.67	125.20	128.60
1	AA	41	A	C5-C6-N6	-5.67	119.17	123.70
51	B3	55	A	C5-C6-N6	-5.67	119.17	123.70
53	B5	720	A	O4'-C1'-N9	5.66	112.73	108.20
53	B5	1552	G	O4'-C1'-N9	5.66	112.73	108.20
53	B5	1610	G	O4'-C1'-N9	5.66	112.73	108.20
53	B5	1940	G	C5-C6-O6	-5.66	125.20	128.60
1	AA	935	G	O4'-C1'-N9	5.66	112.73	108.20
1	AA	523	G	C5-C6-O6	-5.66	125.20	128.60
52	B4	107	G	C5-C6-O6	-5.66	125.20	128.60
53	B5	2258	U	O4'-C1'-N1	5.66	112.73	108.20
1	AA	755	A	O4'-C1'-N9	5.66	112.73	108.20
52	B4	80	A	O4'-C1'-N9	5.66	112.73	108.20
53	B5	1005	G	C5-C6-O6	-5.66	125.20	128.60
53	B5	2313	A	O4'-C1'-N9	5.66	112.73	108.20
1	AA	1643	G	O4'-C1'-N9	5.66	112.73	108.20
53	B5	6	A	O4'-C1'-N9	5.66	112.73	108.20
1	AA	363	G	O4'-C1'-N9	5.66	112.72	108.20
1	AA	1045	G	O4'-C1'-N9	5.66	112.72	108.20
1	AA	1111	G	O4'-C1'-N9	5.66	112.72	108.20
17	AR	284	ALA	C-N-CA	-5.66	107.56	121.70
53	B5	965	A	C5-C6-N6	-5.66	119.17	123.70
53	B5	600	G	O4'-C1'-N9	5.65	112.72	108.20
53	B5	1087	G	N1-C6-O6	5.65	123.29	119.90
53	B5	2964	G	O4'-C1'-N9	5.65	112.72	108.20
1	AA	1162	G	C5-C6-O6	-5.65	125.21	128.60
1	AA	1261	G	C5-C6-O6	-5.65	125.21	128.60
1	AA	1387	U	O4'-C1'-N1	5.65	112.72	108.20
53	B5	136	G	C5-C6-O6	-5.65	125.21	128.60
53	B5	1739	U	O4'-C1'-N1	5.65	112.72	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	2226	U	O4'-C1'-N1	5.65	112.72	108.20
53	B5	2387	A	C5-C6-N6	-5.65	119.18	123.70
53	B5	2480	A	C5-C6-N6	-5.65	119.18	123.70
53	B5	3330	A	O4'-C1'-N9	5.65	112.72	108.20
1	AA	210	A	O4'-C1'-N9	5.65	112.72	108.20
1	AA	403	G	C5-C6-O6	-5.65	125.21	128.60
1	AA	1269	G	N1-C6-O6	5.65	123.29	119.90
1	AA	1348	G	N1-C6-O6	5.65	123.29	119.90
1	AA	1413	G	N1-C6-O6	5.65	123.29	119.90
53	B5	604	G	O4'-C1'-N9	5.65	112.72	108.20
53	B5	1413	G	O4'-C1'-N9	5.65	112.72	108.20
53	B5	2409	G	C5-C6-O6	-5.65	125.21	128.60
1	AA	1273	G	O4'-C1'-N9	5.65	112.72	108.20
1	AA	1308	U	O4'-C1'-N1	5.65	112.72	108.20
53	B5	2698	G	O4'-C1'-N9	5.65	112.72	108.20
53	B5	334	A	O4'-C1'-N9	5.65	112.72	108.20
53	B5	657	A	O4'-C1'-N9	5.65	112.72	108.20
53	B5	845	G	C5-C6-O6	-5.65	125.21	128.60
53	B5	2386	A	C5-C6-N6	-5.65	119.18	123.70
1	AA	869	A	O4'-C1'-N9	5.65	112.72	108.20
29	BE	207	TYR	CB-CG-CD2	-5.65	117.61	121.00
53	B5	2805	G	C5-C6-O6	-5.65	125.21	128.60
1	AA	456	A	C5-C6-N6	-5.64	119.18	123.70
1	AA	494	U	O4'-C1'-N1	5.64	112.72	108.20
53	B5	3	U	O4'-C1'-N1	5.64	112.72	108.20
19	A7	1	G	C8-N9-C4	5.64	108.66	106.40
53	B5	1679	A	C5-C6-N6	-5.64	119.19	123.70
53	B5	2503	G	C5-C6-O6	-5.64	125.22	128.60
53	B5	388	G	N1-C6-O6	5.64	123.28	119.90
1	AA	39	A	O4'-C1'-N9	5.64	112.71	108.20
1	AA	509	G	O4'-C1'-N9	5.64	112.71	108.20
1	AA	510	G	O4'-C1'-N9	5.64	112.71	108.20
1	AA	732	G	O4'-C1'-N9	5.64	112.71	108.20
1	AA	899	A	O4'-C1'-N9	5.64	112.71	108.20
1	AA	1502	G	O4'-C1'-N9	5.64	112.71	108.20
53	B5	1147	G	N1-C6-O6	5.64	123.28	119.90
53	B5	1428	A	O4'-C1'-N9	5.64	112.71	108.20
53	B5	3129	A	O4'-C1'-N9	5.64	112.71	108.20
1	AA	1258	U	O4'-C1'-N1	5.64	112.71	108.20
53	B5	236	G	O4'-C1'-N9	5.64	112.71	108.20
53	B5	717	C	C6-N1-C1'	-5.64	114.03	120.80
1	AA	1710	A	O4'-C1'-N9	5.64	112.71	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	91	G	C5-C6-O6	-5.64	125.22	128.60
53	B5	668	G	C5-C6-O6	-5.64	125.22	128.60
53	B5	817	A	C5-C6-N6	-5.64	119.19	123.70
53	B5	976	C	N3-C4-C5	-5.64	119.65	121.90
53	B5	1573	G	O4'-C1'-N9	5.64	112.71	108.20
53	B5	1744	G	O4'-C1'-N9	5.64	112.71	108.20
53	B5	2381	G	O4'-C1'-N9	5.64	112.71	108.20
53	B5	683	U	O4'-C1'-N1	5.63	112.71	108.20
1	AA	444	C	O4'-C1'-N1	5.63	112.71	108.20
1	AA	1548	A	C5'-C4'-C3'	-5.63	106.99	116.00
1	AA	1743	G	C5-C6-O6	-5.63	125.22	128.60
51	B3	114	A	O4'-C1'-N9	5.63	112.71	108.20
53	B5	2939	G	N1-C6-O6	5.63	123.28	119.90
1	AA	42	G	P-O3'-C3'	5.63	126.46	119.70
53	B5	3361	G	C5-C6-O6	-5.63	125.22	128.60
53	B5	1289	G	O4'-C1'-N9	5.63	112.70	108.20
1	AA	68	A	O4'-C1'-N9	5.63	112.70	108.20
1	AA	803	A	O4'-C1'-N9	5.63	112.70	108.20
51	B3	69	G	N1-C6-O6	5.63	123.28	119.90
53	B5	1249	G	N1-C6-O6	5.63	123.28	119.90
53	B5	1313	G	C5'-C4'-O4'	5.63	115.85	109.10
1	AA	105	A	O4'-C1'-N9	5.62	112.70	108.20
1	AA	127	G	O4'-C1'-N9	5.62	112.70	108.20
1	AA	1266	G	N1-C6-O6	5.62	123.27	119.90
53	B5	264	G	C5-C6-O6	-5.62	125.23	128.60
53	B5	774	G	O4'-C1'-N9	5.62	112.70	108.20
53	B5	997	A	C5-C6-N6	-5.62	119.20	123.70
53	B5	1273	A	O4'-C1'-N9	5.62	112.70	108.20
53	B5	1561	G	O4'-C1'-N9	5.62	112.70	108.20
53	B5	2232	A	O4'-C1'-N9	5.62	112.70	108.20
53	B5	2609	A	C5-C6-N6	-5.62	119.20	123.70
53	B5	3256	G	O4'-C1'-N9	5.62	112.70	108.20
53	B5	3368	U	O4'-C1'-N1	5.62	112.70	108.20
1	AA	451	A	P-O3'-C3'	5.62	126.44	119.70
1	AA	1510	G	O4'-C1'-N9	5.62	112.70	108.20
53	B5	330	G	C5-C6-O6	-5.62	125.23	128.60
53	B5	714	G	P-O3'-C3'	5.62	126.45	119.70
53	B5	1245	A	C5-C6-N6	-5.62	119.20	123.70
53	B5	1300	G	C5-C6-O6	-5.62	125.23	128.60
53	B5	1456	A	C5-C6-N6	-5.62	119.20	123.70
53	B5	1654	A	C5'-C4'-O4'	5.62	115.84	109.10
1	AA	871	G	O4'-C1'-N9	5.62	112.70	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	956	G	O4'-C1'-N9	5.62	112.70	108.20
1	AA	1511	G	O4'-C1'-N9	5.62	112.70	108.20
19	A7	62	A	O4'-C1'-C2'	5.62	112.66	107.60
53	B5	3027	A	P-O3'-C3'	5.62	126.44	119.70
53	B5	3230	G	C5-C6-O6	-5.62	125.23	128.60
1	AA	1234	A	C5-C6-N6	-5.62	119.20	123.70
53	B5	913	A	O4'-C1'-N9	5.62	112.69	108.20
53	B5	1602	A	C5-C6-N6	-5.62	119.20	123.70
53	B5	2460	U	O4'-C1'-N1	5.62	112.69	108.20
53	B5	2671	A	C5-C6-N6	-5.62	119.21	123.70
1	AA	235	G	O4'-C1'-N9	5.62	112.69	108.20
1	AA	853	G	C5-C6-O6	-5.62	125.23	128.60
1	AA	1647	G	C5-C6-O6	-5.62	125.23	128.60
53	B5	412	G	C5-C6-O6	-5.62	125.23	128.60
53	B5	473	G	C5-C6-O6	-5.62	125.23	128.60
53	B5	625	G	O4'-C1'-N9	5.62	112.69	108.20
1	AA	922	A	O4'-C1'-N9	5.61	112.69	108.20
53	B5	1488	G	O4'-C1'-N9	5.61	112.69	108.20
53	B5	1646	G	C5-C6-O6	-5.61	125.23	128.60
53	B5	2137	U	O4'-C1'-N1	5.61	112.69	108.20
53	B5	3052	G	O4'-C1'-N9	5.61	112.69	108.20
53	B5	3263	G	O4'-C1'-N9	5.61	112.69	108.20
19	A7	23	A	C5-C6-N1	5.61	120.50	117.70
53	B5	3119	U	O4'-C1'-N1	5.61	112.69	108.20
1	AA	1168	A	O4'-C1'-N9	5.61	112.69	108.20
1	AA	244	A	O4'-C1'-N9	5.61	112.69	108.20
1	AA	995	U	O4'-C1'-N1	5.61	112.69	108.20
1	AA	1784	G	C5-C6-O6	-5.61	125.24	128.60
53	B5	161	G	N1-C6-O6	5.61	123.26	119.90
53	B5	832	G	C5-C6-O6	-5.61	125.23	128.60
53	B5	902	G	O4'-C1'-N9	5.61	112.69	108.20
53	B5	2288	G	C5-C6-O6	-5.61	125.23	128.60
53	B5	2538	U	O4'-C1'-N1	5.61	112.69	108.20
53	B5	2661	G	O4'-C1'-N9	5.61	112.69	108.20
1	AA	28	A	O4'-C1'-N9	5.61	112.68	108.20
1	AA	273	G	C5-C6-O6	-5.61	125.24	128.60
1	AA	345	U	O4'-C1'-N1	5.61	112.69	108.20
1	AA	355	G	N1-C6-O6	5.61	123.26	119.90
1	AA	1107	G	O4'-C1'-N9	5.61	112.68	108.20
1	AA	1225	G	N1-C6-O6	5.61	123.26	119.90
1	AA	1521	G	N1-C6-O6	-5.61	116.54	119.90
53	B5	1212	A	O4'-C1'-N9	5.61	112.68	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	2615	G	C5-C6-O6	-5.61	125.24	128.60
1	AA	829	A	O4'-C1'-N9	5.60	112.68	108.20
1	AA	1697	G	O4'-C1'-N9	5.60	112.68	108.20
53	B5	136	G	O4'-C1'-N9	5.60	112.68	108.20
53	B5	2557	G	N1-C6-O6	5.60	123.26	119.90
1	AA	197	A	O4'-C1'-N9	5.60	112.68	108.20
1	AA	329	G	O4'-C1'-N9	5.60	112.68	108.20
1	AA	1119	G	C5-C6-O6	-5.60	125.24	128.60
53	B5	847	A	C5-C6-N6	-5.60	119.22	123.70
53	B5	1027	A	P-O3'-C3'	5.60	126.42	119.70
53	B5	2447	A	C5-C6-N6	-5.60	119.22	123.70
1	AA	1129	A	O4'-C1'-N9	5.60	112.68	108.20
19	A7	24	G	C8-N9-C4	-5.60	104.16	106.40
51	B3	78	G	C5-C6-O6	-5.60	125.24	128.60
53	B5	170	G	O4'-C1'-N9	5.60	112.68	108.20
53	B5	1311	G	O4'-C1'-N9	5.60	112.68	108.20
1	AA	770	A	O4'-C1'-N9	5.60	112.68	108.20
1	AA	1119	G	O4'-C1'-N9	5.60	112.68	108.20
1	AA	1196	G	O4'-C4'-C3'	-5.60	98.40	104.00
53	B5	287	G	C5-C6-O6	-5.60	125.24	128.60
53	B5	477	A	O4'-C1'-N9	5.60	112.68	108.20
53	B5	1207	G	C5-C6-O6	-5.60	125.24	128.60
53	B5	2700	G	C5-C6-O6	-5.60	125.24	128.60
1	AA	538	A	O4'-C1'-N9	5.60	112.68	108.20
1	AA	1115	G	O4'-C1'-N9	5.60	112.68	108.20
1	AA	1143	G	C5-C6-O6	-5.60	125.24	128.60
1	AA	1392	G	O4'-C1'-N9	5.60	112.68	108.20
1	AA	1631	A	OP1-P-OP2	-5.60	111.20	119.60
52	B4	46	G	C5-C6-O6	-5.60	125.24	128.60
53	B5	810	A	C4-C5-C6	5.60	119.80	117.00
53	B5	2289	U	O4'-C1'-N1	5.60	112.68	108.20
53	B5	2375	G	O4'-C1'-N9	5.60	112.68	108.20
53	B5	2730	G	C5-C6-O6	-5.60	125.24	128.60
53	B5	2868	U	O4'-C1'-N1	5.60	112.68	108.20
1	AA	322	G	N1-C6-O6	5.60	123.26	119.90
53	B5	104	G	C5-C6-O6	-5.60	125.24	128.60
53	B5	400	G	O4'-C1'-N9	5.60	112.68	108.20
53	B5	887	G	O4'-C1'-N9	5.60	112.68	108.20
53	B5	3041	U	O4'-C1'-N1	5.60	112.68	108.20
1	AA	689	G	O4'-C1'-N9	5.59	112.67	108.20
1	AA	859	A	C8-N9-C4	-5.59	103.56	105.80
53	B5	3016	A	C4-C5-C6	5.59	119.80	117.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	3146	G	C1'-O4'-C4'	-5.59	105.42	109.90
1	AA	845	G	O4'-C1'-N9	5.59	112.67	108.20
53	B5	133	U	O4'-C1'-N1	5.59	112.67	108.20
53	B5	2311	G	O4'-C1'-N9	5.59	112.67	108.20
19	A7	56	C	N1-C2-O2	5.59	122.25	118.90
53	B5	428	A	O4'-C1'-N9	5.59	112.67	108.20
53	B5	646	A	O4'-C1'-N9	5.59	112.67	108.20
53	B5	1213	G	O4'-C1'-N9	5.59	112.67	108.20
53	B5	2846	U	O4'-C1'-N1	5.59	112.67	108.20
53	B5	2957	G	O4'-C1'-N9	5.59	112.67	108.20
1	AA	1091	G	O4'-C1'-N9	5.59	112.67	108.20
53	B5	1099	A	O4'-C1'-N9	5.59	112.67	108.20
53	B5	1147	G	O4'-C1'-N9	5.59	112.67	108.20
53	B5	2488	A	C4-C5-C6	5.59	119.80	117.00
53	B5	2585	G	C5-C6-O6	-5.59	125.25	128.60
53	B5	2364	G	C5-C6-O6	-5.59	125.25	128.60
53	B5	2632	G	O4'-C1'-N9	5.59	112.67	108.20
53	B5	2749	G	O4'-C1'-N9	5.59	112.67	108.20
53	B5	2800	G	O4'-C1'-N9	5.59	112.67	108.20
1	AA	288	A	O4'-C1'-N9	5.59	112.67	108.20
1	AA	997	A	O4'-C1'-N9	5.59	112.67	108.20
1	AA	1364	G	N1-C6-O6	5.59	123.25	119.90
53	B5	387	A	O4'-C1'-N9	5.59	112.67	108.20
53	B5	1542	G	C5-C6-O6	-5.59	125.25	128.60
1	AA	259	U	O4'-C1'-N1	5.58	112.67	108.20
53	B5	1097	G	O4'-C1'-N9	5.58	112.67	108.20
53	B5	1237	G	N1-C6-O6	5.58	123.25	119.90
53	B5	1929	G	O4'-C1'-N9	5.58	112.67	108.20
53	B5	2818	U	O4'-C1'-N1	5.58	112.67	108.20
1	AA	390	G	N1-C6-O6	5.58	123.25	119.90
1	AA	419	G	C5-C6-O6	-5.58	125.25	128.60
1	AA	1162	G	O4'-C1'-N9	5.58	112.67	108.20
51	B3	75	G	O4'-C1'-N9	5.58	112.67	108.20
53	B5	341	G	O4'-C1'-N9	5.58	112.67	108.20
53	B5	628	A	C5-C6-N6	-5.58	119.23	123.70
53	B5	1446	A	O4'-C1'-N9	5.58	112.67	108.20
1	AA	898	G	O4'-C1'-N9	5.58	112.67	108.20
1	AA	947	G	O4'-C1'-N9	5.58	112.67	108.20
1	AA	1392	G	C5-C6-O6	-5.58	125.25	128.60
53	B5	16	A	P-O3'-C3'	5.58	126.40	119.70
53	B5	265	A	O4'-C1'-N9	5.58	112.66	108.20
53	B5	725	G	O4'-C1'-N9	5.58	112.67	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	2160	G	O4'-C1'-N9	5.58	112.67	108.20
53	B5	3322	A	C5-C6-N6	-5.58	119.23	123.70
53	B5	145	G	C5-C6-O6	-5.58	125.25	128.60
1	AA	673	A	O4'-C1'-N9	5.58	112.66	108.20
1	AA	729	G	N1-C6-O6	5.58	123.25	119.90
53	B5	712	G	C4-N9-C1'	5.58	133.75	126.50
53	B5	1488	G	C5-C6-O6	-5.58	125.25	128.60
53	B5	2770	G	P-O3'-C3'	5.58	126.39	119.70
53	B5	902	G	N1-C6-O6	5.58	123.25	119.90
53	B5	1250	G	C5-C6-O6	-5.58	125.25	128.60
53	B5	2386	A	O4'-C1'-N9	5.58	112.66	108.20
1	AA	194	U	C2-N1-C1'	5.57	124.39	117.70
53	B5	33	G	O4'-C1'-N9	5.57	112.66	108.20
53	B5	206	G	O4'-C1'-N9	5.57	112.66	108.20
53	B5	1051	U	O4'-C1'-N1	5.57	112.66	108.20
53	B5	1083	G	C5-C6-O6	-5.57	125.26	128.60
53	B5	3102	G	O4'-C1'-N9	5.57	112.66	108.20
52	B4	19	C	N3-C4-N4	5.57	121.90	118.00
53	B5	1480	G	C5-C6-O6	-5.57	125.26	128.60
53	B5	2224	A	C5-C6-N6	-5.57	119.24	123.70
1	AA	971	G	O4'-C1'-N9	5.57	112.66	108.20
1	AA	1715	G	O4'-C1'-N9	5.57	112.66	108.20
1	AA	1744	A	O4'-C1'-N9	5.57	112.66	108.20
53	B5	406	G	C5-C6-O6	-5.57	125.26	128.60
53	B5	1190	A	P-O3'-C3'	5.57	126.38	119.70
53	B5	1375	G	O4'-C1'-N9	5.57	112.66	108.20
53	B5	2560	U	O4'-C1'-N1	5.57	112.66	108.20
53	B5	2648	G	C5-C6-O6	-5.57	125.26	128.60
1	AA	300	A	O4'-C1'-N9	5.57	112.66	108.20
1	AA	438	A	O4'-C1'-N9	5.57	112.66	108.20
1	AA	982	A	O4'-C1'-N9	5.57	112.66	108.20
1	AA	996	G	C5-C6-O6	-5.57	125.26	128.60
19	A7	42	G	N9-C4-C5	5.57	107.63	105.40
53	B5	537	A	O4'-C1'-N9	5.57	112.66	108.20
53	B5	1252	A	O4'-C1'-N9	5.57	112.66	108.20
53	B5	1812	G	N1-C6-O6	5.57	123.24	119.90
53	B5	1888	U	O4'-C1'-N1	5.57	112.66	108.20
53	B5	2440	G	C5-C6-O6	-5.57	125.26	128.60
1	AA	763	G	O4'-C1'-N9	5.57	112.65	108.20
1	AA	1068	U	O4'-C1'-N1	5.57	112.65	108.20
1	AA	1660	G	C5-C6-O6	-5.57	125.26	128.60
53	B5	277	G	C5-C6-O6	-5.57	125.26	128.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	1385	C	N3-C4-C5	-5.57	119.67	121.90
53	B5	1637	A	O4'-C1'-N9	5.57	112.65	108.20
1	AA	1741	U	O4'-C1'-N1	5.57	112.65	108.20
18	AT	33	TYR	CB-CG-CD1	5.57	124.34	121.00
51	B3	25	G	C5-C6-O6	-5.57	125.26	128.60
53	B5	1680	G	C5-C6-O6	-5.57	125.26	128.60
53	B5	1885	U	O4'-C1'-N1	5.57	112.65	108.20
53	B5	2634	U	O4'-C1'-N1	5.57	112.65	108.20
53	B5	2659	G	O4'-C1'-N9	5.57	112.65	108.20
53	B5	3078	U	O4'-C1'-N1	5.57	112.65	108.20
1	AA	865	A	OP1-P-OP2	5.56	127.95	119.60
53	B5	33	G	C5-C6-O6	-5.56	125.26	128.60
53	B5	900	G	O4'-C1'-N9	5.56	112.65	108.20
53	B5	2149	A	O4'-C1'-N9	5.56	112.65	108.20
53	B5	3069	G	O4'-C1'-N9	5.56	112.65	108.20
1	AA	642	G	O4'-C1'-N9	5.56	112.65	108.20
1	AA	1003	U	O4'-C1'-N1	5.56	112.65	108.20
53	B5	337	G	C5-C6-O6	-5.56	125.26	128.60
53	B5	835	G	C5-C6-O6	-5.56	125.26	128.60
53	B5	1528	G	C5-C6-O6	-5.56	125.26	128.60
53	B5	1575	A	O4'-C1'-N9	5.56	112.65	108.20
53	B5	659	G	C5-C6-O6	-5.56	125.26	128.60
53	B5	1571	A	O4'-C1'-N9	5.56	112.65	108.20
1	AA	6	G	C5-C6-O6	-5.56	125.26	128.60
1	AA	811	A	O4'-C1'-N9	5.56	112.65	108.20
19	A7	24	G	C4-C5-N7	5.56	113.02	110.80
53	B5	1271	A	C4-C5-C6	5.56	119.78	117.00
53	B5	2550	U	O4'-C1'-N1	5.56	112.65	108.20
1	AA	562	G	O4'-C1'-N9	5.56	112.65	108.20
1	AA	1039	G	C5-C6-O6	-5.56	125.27	128.60
1	AA	1179	A	O4'-C1'-N9	5.56	112.64	108.20
1	AA	1640	G	O4'-C1'-N9	5.56	112.65	108.20
51	B3	21	G	O4'-C1'-N9	5.56	112.65	108.20
53	B5	432	G	C5-C6-O6	-5.56	125.27	128.60
53	B5	2533	G	O4'-C1'-N9	5.56	112.65	108.20
53	B5	3124	G	C5-C6-O6	-5.56	125.27	128.60
53	B5	1789	G	C5-C6-O6	-5.56	125.27	128.60
53	B5	3266	G	N1-C6-O6	5.56	123.23	119.90
19	A7	67	A	C3'-C2'-C1'	5.55	105.94	101.50
53	B5	1	G	N1-C6-O6	5.55	123.23	119.90
53	B5	2395	G	O4'-C1'-N9	5.55	112.64	108.20
53	B5	2451	G	C5-C6-O6	-5.55	125.27	128.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	2747	A	O4'-C1'-N9	5.55	112.64	108.20
1	AA	1140	A	O4'-C1'-N9	5.55	112.64	108.20
1	AA	1194	G	P-O3'-C3'	5.55	126.36	119.70
53	B5	1849	C	O4'-C1'-N1	5.55	112.64	108.20
1	AA	821	U	O4'-C1'-N1	5.55	112.64	108.20
1	AA	895	G	O4'-C1'-N9	5.55	112.64	108.20
1	AA	1080	G	O4'-C1'-N9	5.55	112.64	108.20
1	AA	1491	A	C5-C6-N6	-5.55	119.26	123.70
1	AA	1772	G	C5-C6-O6	-5.55	125.27	128.60
51	B3	94	A	O4'-C1'-N9	5.55	112.64	108.20
53	B5	738	A	O4'-C1'-N9	5.55	112.64	108.20
53	B5	2779	A	O4'-C1'-N9	5.55	112.64	108.20
53	B5	70	A	C4-C5-C6	5.55	119.77	117.00
53	B5	1057	A	O4'-C1'-N9	5.55	112.64	108.20
53	B5	2221	G	O4'-C1'-N9	5.55	112.64	108.20
53	B5	2281	A	O4'-C1'-N9	5.55	112.64	108.20
53	B5	2754	G	C5-C6-O6	-5.55	125.27	128.60
53	B5	2814	G	O4'-C1'-N9	5.55	112.64	108.20
53	B5	2896	A	O4'-C1'-N9	5.55	112.64	108.20
1	AA	723	G	C5-C6-O6	-5.55	125.27	128.60
1	AA	1518	U	O5'-P-OP2	5.55	117.36	110.70
53	B5	1673	G	P-O3'-C3'	5.55	126.36	119.70
53	B5	1727	G	O4'-C1'-N9	5.55	112.64	108.20
53	B5	3016	A	C5-C6-N6	-5.55	119.26	123.70
1	AA	1554	A	O4'-C1'-N9	5.55	112.64	108.20
53	B5	718	G	O4'-C1'-N9	5.55	112.64	108.20
53	B5	1236	G	N1-C6-O6	5.55	123.23	119.90
53	B5	1491	A	C5-C6-N6	-5.55	119.26	123.70
53	B5	2412	G	C5-C6-O6	-5.55	125.27	128.60
53	B5	2670	G	N1-C6-O6	5.55	123.23	119.90
53	B5	2780	A	O4'-C1'-N9	5.55	112.64	108.20
53	B5	2341	A	O4'-C1'-N9	5.54	112.64	108.20
53	B5	3252	G	O4'-C1'-N9	5.54	112.64	108.20
1	AA	1647	G	O4'-C1'-N9	5.54	112.64	108.20
19	A7	53	G	C4'-C3'-C2'	-5.54	97.06	102.60
53	B5	86	G	C5-C6-O6	-5.54	125.27	128.60
53	B5	866	A	O4'-C1'-N9	5.54	112.64	108.20
53	B5	870	G	O4'-C1'-N9	5.54	112.64	108.20
53	B5	1064	A	O4'-C1'-N9	5.54	112.64	108.20
53	B5	1798	A	O4'-C1'-N9	5.54	112.64	108.20
53	B5	2811	A	C5-C6-N6	-5.54	119.27	123.70
53	B5	3218	A	C5-C6-N6	-5.54	119.27	123.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	328	A	O4'-C1'-N9	5.54	112.63	108.20
1	AA	1220	A	O4'-C1'-N9	5.54	112.63	108.20
17	AR	286	GLU	CA-CB-CG	5.54	125.59	113.40
19	A7	68	U	N1-C2-O2	5.54	126.68	122.80
53	B5	13	A	C5-C6-N6	-5.54	119.27	123.70
53	B5	39	A	C4-C5-C6	5.54	119.77	117.00
53	B5	582	G	C5-C6-O6	-5.54	125.28	128.60
53	B5	1744	G	C5-C6-O6	-5.54	125.28	128.60
1	AA	154	G	C5-C6-O6	-5.54	125.28	128.60
1	AA	170	U	O4'-C1'-N1	5.54	112.63	108.20
53	B5	236	G	C5-C6-O6	-5.54	125.28	128.60
1	AA	989	C	C1'-O4'-C4'	-5.54	105.47	109.90
53	B5	979	G	C5-C6-O6	-5.54	125.28	128.60
53	B5	1729	A	O4'-C1'-N9	5.54	112.63	108.20
53	B5	1775	G	O4'-C1'-N9	5.54	112.63	108.20
53	B5	3337	G	C5-C6-O6	-5.54	125.28	128.60
1	AA	1729	A	O4'-C1'-N9	5.54	112.63	108.20
53	B5	160	G	C5-C6-O6	-5.54	125.28	128.60
1	AA	247	A	O4'-C1'-N9	5.54	112.63	108.20
51	B3	41	G	C5-C6-O6	-5.54	125.28	128.60
51	B3	80	G	O4'-C1'-N9	5.54	112.63	108.20
53	B5	243	G	N1-C6-O6	5.54	123.22	119.90
53	B5	1230	G	C5-C6-O6	-5.54	125.28	128.60
53	B5	2665	U	O4'-C1'-N1	5.54	112.63	108.20
53	B5	2728	G	C5-C6-O6	-5.54	125.28	128.60
53	B5	3140	G	C5-C6-O6	-5.54	125.28	128.60
53	B5	3256	G	C5-C6-O6	-5.54	125.28	128.60
1	AA	1301	G	N1-C6-O6	5.53	123.22	119.90
53	B5	1384	U	C2-N1-C1'	5.53	124.34	117.70
53	B5	2235	C	N3-C4-N4	5.53	121.87	118.00
53	B5	2236	G	O4'-C1'-N9	5.53	112.63	108.20
53	B5	2753	G	C5-C6-O6	-5.53	125.28	128.60
1	AA	871	G	C5-C6-O6	-5.53	125.28	128.60
53	B5	324	A	C5-C6-N6	-5.53	119.28	123.70
53	B5	2669	G	O4'-C1'-N9	5.53	112.62	108.20
53	B5	386	A	C5-C6-N6	-5.53	119.28	123.70
53	B5	936	A	O4'-C1'-N9	5.53	112.62	108.20
53	B5	2283	G	C5-C6-O6	-5.53	125.28	128.60
53	B5	2377	G	O4'-C1'-N9	5.53	112.62	108.20
53	B5	2466	G	C5-C6-O6	-5.53	125.28	128.60
53	B5	3103	A	O4'-C1'-N9	5.53	112.62	108.20
53	B5	3239	G	C5-C6-O6	-5.53	125.28	128.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1195	G	C5-C6-O6	-5.53	125.28	128.60
1	AA	866	G	C6-C5-N7	-5.53	127.08	130.40
1	AA	1022	A	C5-C6-N6	-5.53	119.28	123.70
1	AA	1106	G	O4'-C1'-N9	5.53	112.62	108.20
53	B5	378	A	O4'-C1'-N9	5.53	112.62	108.20
53	B5	965	A	O4'-C1'-N9	5.53	112.62	108.20
53	B5	1401	A	O4'-C1'-N9	5.53	112.62	108.20
53	B5	1711	C	N3-C4-N4	5.53	121.87	118.00
53	B5	2279	A	C4-C5-C6	5.53	119.76	117.00
1	AA	180	A	O4'-C1'-N9	5.53	112.62	108.20
1	AA	549	G	O4'-C1'-N9	5.53	112.62	108.20
1	AA	950	A	O4'-C1'-N9	5.53	112.62	108.20
1	AA	992	A	O4'-C1'-N9	5.53	112.62	108.20
53	B5	57	A	O4'-C1'-N9	5.53	112.62	108.20
53	B5	187	A	O4'-C1'-N9	5.53	112.62	108.20
53	B5	1282	G	O4'-C1'-N9	5.53	112.62	108.20
1	AA	1399	G	O4'-C1'-N9	5.52	112.62	108.20
53	B5	2834	G	C5-C6-O6	-5.52	125.29	128.60
1	AA	1190	A	O4'-C1'-N9	5.52	112.62	108.20
1	AA	1292	G	C5-C6-O6	-5.52	125.29	128.60
53	B5	198	A	O4'-C1'-N9	5.52	112.62	108.20
53	B5	1234	G	C5-C6-O6	-5.52	125.29	128.60
53	B5	1623	G	C5-C6-O6	-5.52	125.29	128.60
53	B5	2367	A	O4'-C1'-N9	5.52	112.62	108.20
53	B5	3326	G	N1-C6-O6	5.52	123.21	119.90
1	AA	1545	A	C4'-C3'-C2'	-5.52	97.08	102.60
53	B5	365	A	O4'-C1'-N9	5.52	112.62	108.20
53	B5	1654	A	C4-C5-C6	5.52	119.76	117.00
1	AA	879	G	O4'-C1'-N9	5.52	112.62	108.20
1	AA	891	A	O4'-C1'-N9	5.52	112.62	108.20
53	B5	997	A	O4'-C1'-N9	5.52	112.62	108.20
53	B5	1059	G	O4'-C1'-N9	5.52	112.61	108.20
53	B5	1374	G	C5-C6-O6	-5.52	125.29	128.60
1	AA	116	U	O4'-C1'-N1	5.52	112.61	108.20
1	AA	383	G	O4'-C1'-N9	5.52	112.61	108.20
53	B5	1064	A	C5-C6-N1	-5.52	114.94	117.70
53	B5	1204	A	O4'-C1'-N9	5.52	112.61	108.20
53	B5	1287	A	O4'-C1'-N9	5.52	112.61	108.20
53	B5	1541	G	C5-C6-O6	-5.52	125.29	128.60
53	B5	1587	A	C4-C5-C6	5.52	119.76	117.00
53	B5	1784	G	C5-C6-O6	-5.52	125.29	128.60
1	AA	1658	A	O4'-C1'-N9	5.52	112.61	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
33	BI	104	SER	N-CA-CB	5.52	118.77	110.50
52	B4	131	A	O4'-C1'-N9	5.52	112.61	108.20
53	B5	622	A	O4'-C1'-N9	5.52	112.61	108.20
53	B5	3008	A	O4'-C1'-N9	5.52	112.61	108.20
1	AA	506	A	O4'-C1'-N9	5.51	112.61	108.20
1	AA	1038	A	O4'-C1'-N9	5.51	112.61	108.20
1	AA	1205	C	N3-C4-N4	5.51	121.86	118.00
8	AI	97	VAL	CB-CA-C	5.51	121.88	111.40
52	B4	16	G	C5-C6-O6	-5.51	125.29	128.60
53	B5	521	A	O4'-C1'-N9	5.51	112.61	108.20
53	B5	904	A	O4'-C1'-N9	5.51	112.61	108.20
53	B5	1856	C	O4'-C1'-N1	5.51	112.61	108.20
53	B5	2395	G	C5-C6-O6	-5.51	125.29	128.60
1	AA	69	G	N1-C6-O6	5.51	123.21	119.90
53	B5	514	G	C5-C6-O6	-5.51	125.29	128.60
53	B5	917	A	C5-C6-N6	-5.51	119.29	123.70
53	B5	1084	A	O4'-C1'-N9	5.51	112.61	108.20
53	B5	1232	C	N3-C4-N4	5.51	121.86	118.00
53	B5	1249	G	P-O3'-C3'	5.51	126.31	119.70
1	AA	246	G	C5-C6-O6	-5.51	125.30	128.60
1	AA	746	A	O4'-C1'-N9	5.51	112.61	108.20
1	AA	1264	G	C5-C6-O6	-5.51	125.29	128.60
1	AA	1292	G	O4'-C1'-N9	5.51	112.61	108.20
53	B5	419	G	C5-C6-O6	-5.51	125.29	128.60
53	B5	607	A	O4'-C1'-N9	5.51	112.61	108.20
53	B5	771	A	O4'-C1'-N9	5.51	112.61	108.20
53	B5	1806	A	P-O3'-C3'	5.51	126.31	119.70
53	B5	1922	A	O4'-C1'-N9	5.51	112.61	108.20
53	B5	2290	C	O4'-C1'-N1	5.51	112.61	108.20
53	B5	3187	U	O4'-C1'-N1	5.51	112.61	108.20
1	AA	153	G	O4'-C1'-N9	5.51	112.61	108.20
53	B5	661	G	C5-C6-O6	-5.51	125.30	128.60
1	AA	1246	U	O4'-C1'-N1	5.51	112.61	108.20
19	A7	28	C	N3-C4-N4	-5.51	114.15	118.00
53	B5	608	A	O4'-C1'-N9	5.51	112.60	108.20
53	B5	2951	G	C5-C6-O6	-5.50	125.30	128.60
1	AA	634	G	C5-C6-O6	-5.50	125.30	128.60
1	AA	1196	G	C5-C6-O6	-5.50	125.30	128.60
1	AA	1792	A	C5-C6-N6	-5.50	119.30	123.70
51	B3	70	A	C5-C6-N6	-5.50	119.30	123.70
52	B4	125	U	O4'-C1'-N1	5.50	112.60	108.20
53	B5	1261	G	N1-C6-O6	5.50	123.20	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	1603	A	C5-C6-N6	-5.50	119.30	123.70
53	B5	1789	G	O4'-C1'-N9	5.50	112.60	108.20
53	B5	2172	A	C5-C6-N6	-5.50	119.30	123.70
1	AA	970	A	O4'-C1'-N9	5.50	112.60	108.20
1	AA	1137	G	O4'-C1'-N9	5.50	112.60	108.20
53	B5	813	G	C5-C6-O6	-5.50	125.30	128.60
53	B5	3189	G	O4'-C1'-N9	5.50	112.60	108.20
27	BC	129	ALA	N-CA-CB	5.50	117.80	110.10
53	B5	856	G	O4'-C1'-N9	5.50	112.60	108.20
53	B5	2574	G	N1-C6-O6	5.50	123.20	119.90
1	AA	173	A	O4'-C1'-N9	5.50	112.60	108.20
1	AA	287	G	C5-C6-O6	-5.50	125.30	128.60
53	B5	34	A	C4-C5-C6	5.50	119.75	117.00
53	B5	287	G	O4'-C1'-N9	5.50	112.60	108.20
53	B5	567	G	O4'-C1'-N9	5.50	112.60	108.20
53	B5	856	G	C5-C6-O6	-5.50	125.30	128.60
53	B5	1538	G	C5-C6-O6	-5.50	125.30	128.60
53	B5	2794	G	O4'-C1'-N9	5.50	112.60	108.20
1	AA	838	G	O4'-C1'-N9	5.50	112.60	108.20
19	A7	8	U	C5-C6-N1	-5.50	119.95	122.70
53	B5	66	A	O4'-C1'-N9	5.50	112.60	108.20
53	B5	77	A	C5-C6-N6	-5.50	119.30	123.70
53	B5	171	G	N1-C6-O6	5.50	123.20	119.90
53	B5	299	G	C5-C6-O6	-5.50	125.30	128.60
53	B5	2255	A	O4'-C1'-N9	5.50	112.60	108.20
53	B5	3017	A	O4'-C1'-N9	5.50	112.60	108.20
53	B5	3396	U	O4'-C1'-N1	5.50	112.60	108.20
1	AA	1518	U	C3'-C2'-C1'	-5.50	97.10	101.50
53	B5	2748	A	O4'-C1'-N9	5.50	112.60	108.20
19	A7	53	G	C2-N3-C4	5.49	114.65	111.90
53	B5	529	A	O4'-C1'-N9	5.49	112.59	108.20
1	AA	112	A	O4'-C1'-N9	5.49	112.59	108.20
1	AA	437	A	C5-C6-N6	-5.49	119.31	123.70
1	AA	925	A	O4'-C1'-N9	5.49	112.59	108.20
53	B5	3317	U	O4'-C1'-N1	5.49	112.59	108.20
1	AA	1435	G	C5-C6-O6	-5.49	125.31	128.60
53	B5	2623	G	O4'-C1'-N9	5.49	112.59	108.20
1	AA	457	G	O4'-C1'-N9	5.49	112.59	108.20
1	AA	601	A	O4'-C1'-N9	5.49	112.59	108.20
1	AA	1062	A	O4'-C1'-N9	5.49	112.59	108.20
1	AA	1415	G	P-O3'-C3'	5.49	126.28	119.70
1	AA	1567	A	O4'-C1'-N9	5.49	112.59	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	214	G	O4'-C1'-N9	5.49	112.59	108.20
53	B5	1418	A	C4-C5-C6	5.49	119.74	117.00
53	B5	3059	G	C5-C6-O6	-5.49	125.31	128.60
53	B5	3282	U	O4'-C1'-N1	5.49	112.59	108.20
52	B4	124	G	C5-C6-O6	-5.49	125.31	128.60
53	B5	188	U	O4'-C1'-N1	5.49	112.59	108.20
53	B5	689	U	O4'-C1'-N1	5.49	112.59	108.20
1	AA	429	G	O4'-C1'-N9	5.49	112.59	108.20
1	AA	432	G	O4'-C1'-N9	5.49	112.59	108.20
1	AA	1216	C	N3-C4-N4	5.49	121.84	118.00
52	B4	46	G	O4'-C1'-N9	5.49	112.59	108.20
53	B5	433	A	C5-C6-N6	-5.49	119.31	123.70
53	B5	517	G	O4'-C1'-N9	5.49	112.59	108.20
53	B5	960	U	C2-N1-C1'	5.49	124.28	117.70
53	B5	1139	G	O4'-C1'-N9	5.49	112.59	108.20
53	B5	2291	A	C5-C6-N1	-5.49	114.96	117.70
53	B5	2295	A	O4'-C1'-N9	5.49	112.59	108.20
53	B5	3322	A	O4'-C1'-N9	5.49	112.59	108.20
53	B5	838	G	O4'-C1'-N9	5.48	112.59	108.20
53	B5	2370	G	O4'-C1'-N9	5.48	112.59	108.20
1	AA	253	A	O4'-C1'-N9	5.48	112.59	108.20
1	AA	1765	G	N1-C6-O6	5.48	123.19	119.90
52	B4	42	G	O4'-C1'-N9	5.48	112.59	108.20
53	B5	279	U	O4'-C1'-N1	5.48	112.59	108.20
53	B5	987	U	O4'-C1'-N1	5.48	112.59	108.20
53	B5	1310	G	O4'-C1'-N9	5.48	112.59	108.20
53	B5	1610	G	C5-C6-O6	-5.48	125.31	128.60
53	B5	3074	G	C5-C6-O6	-5.48	125.31	128.60
53	B5	3305	A	C4-C5-C6	5.48	119.74	117.00
1	AA	39	A	C5-C6-N6	-5.48	119.31	123.70
1	AA	156	A	O4'-C1'-N9	5.48	112.58	108.20
1	AA	837	G	O4'-C1'-N9	5.48	112.58	108.20
1	AA	1116	G	C5-C6-O6	-5.48	125.31	128.60
1	AA	1407	A	O4'-C1'-N9	5.48	112.58	108.20
53	B5	616	G	O4'-C1'-N9	5.48	112.58	108.20
53	B5	875	G	C5-C6-O6	-5.48	125.31	128.60
53	B5	3206	A	O4'-C1'-N9	5.48	112.58	108.20
1	AA	985	G	C5-C6-O6	-5.48	125.31	128.60
53	B5	2718	U	C5'-C4'-O4'	5.48	115.67	109.10
1	AA	10	G	O4'-C1'-N9	5.48	112.58	108.20
1	AA	1610	U	O4'-C1'-N1	5.48	112.58	108.20
1	AA	1668	G	O4'-C1'-N9	5.48	112.58	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	781	G	C5-C6-O6	-5.48	125.31	128.60
53	B5	891	G	O4'-C1'-N9	5.48	112.58	108.20
53	B5	1812	G	O4'-C1'-N9	5.48	112.58	108.20
53	B5	2294	U	O4'-C1'-N1	5.48	112.58	108.20
53	B5	2447	A	O4'-C1'-N9	5.48	112.58	108.20
53	B5	134	U	O4'-C1'-N1	5.48	112.58	108.20
53	B5	2169	G	C5-C6-O6	-5.48	125.31	128.60
53	B5	2326	A	C5-C6-N6	-5.48	119.32	123.70
53	B5	2984	C	N3-C4-N4	5.48	121.83	118.00
1	AA	195	G	O4'-C1'-N9	5.47	112.58	108.20
1	AA	264	G	N1-C6-O6	5.47	123.18	119.90
1	AA	383	G	C5-C6-O6	-5.47	125.31	128.60
53	B5	127	G	O4'-C1'-N9	5.47	112.58	108.20
53	B5	1207	G	O4'-C1'-N9	5.47	112.58	108.20
53	B5	2249	G	N1-C6-O6	5.47	123.19	119.90
53	B5	256	G	O4'-C1'-N9	5.47	112.58	108.20
1	AA	260	U	O4'-C1'-N1	5.47	112.58	108.20
1	AA	1234	A	O4'-C1'-N9	5.47	112.58	108.20
53	B5	3343	G	C5-C6-O6	-5.47	125.32	128.60
53	B5	3367	C	P-O3'-C3'	5.47	126.27	119.70
1	AA	650	U	O4'-C1'-N1	5.47	112.58	108.20
1	AA	978	A	O4'-C1'-N9	5.47	112.58	108.20
1	AA	1305	G	C5-C6-O6	-5.47	125.32	128.60
1	AA	1476	G	O4'-C1'-N9	5.47	112.58	108.20
28	BD	50	TYR	CB-CG-CD2	5.47	124.28	121.00
53	B5	1643	A	C5-C6-N6	-5.47	119.33	123.70
1	AA	437	A	O4'-C1'-N9	5.47	112.58	108.20
53	B5	65	A	P-O3'-C3'	5.47	126.26	119.70
1	AA	364	G	O4'-C1'-N9	5.47	112.57	108.20
53	B5	1382	G	C5-C6-O6	-5.47	125.32	128.60
53	B5	1497	C	N3-C4-N4	5.47	121.83	118.00
53	B5	1721	U	O4'-C1'-N1	5.47	112.57	108.20
53	B5	2385	G	O4'-C1'-N9	5.47	112.57	108.20
53	B5	2453	U	O4'-C1'-N1	5.47	112.57	108.20
53	B5	2847	A	C5-C6-N6	-5.47	119.33	123.70
1	AA	84	A	O4'-C1'-N9	5.46	112.57	108.20
1	AA	866	G	O5'-C5'-C4'	5.46	122.08	111.70
51	B3	99	A	O4'-C1'-N9	5.46	112.57	108.20
52	B4	18	U	O4'-C1'-N1	5.46	112.57	108.20
53	B5	635	G	C5-C6-O6	-5.46	125.32	128.60
53	B5	1090	G	C5-C6-O6	-5.46	125.32	128.60
53	B5	3061	G	O4'-C1'-N9	5.46	112.57	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	422	G	O4'-C1'-N9	5.46	112.57	108.20
53	B5	2987	A	O4'-C1'-N9	5.46	112.57	108.20
1	AA	1172	U	O4'-C1'-N1	5.46	112.57	108.20
1	AA	1634	C	C1'-O4'-C4'	-5.46	105.53	109.90
19	A7	21	A	P-O3'-C3'	5.46	126.25	119.70
19	A7	43	G	C5-C6-O6	5.46	131.88	128.60
53	B5	137	G	C5-C6-O6	-5.46	125.32	128.60
53	B5	490	A	O4'-C1'-N9	5.46	112.57	108.20
53	B5	869	G	O4'-C1'-N9	5.46	112.57	108.20
53	B5	891	G	C5-C6-O6	-5.46	125.32	128.60
53	B5	2115	G	C5-C6-O6	-5.46	125.32	128.60
53	B5	2201	G	C5-C6-O6	-5.46	125.32	128.60
53	B5	2368	A	O4'-C1'-N9	5.46	112.57	108.20
53	B5	2703	A	C5-C6-N6	-5.46	119.33	123.70
53	B5	3323	A	O4'-C1'-N9	5.46	112.57	108.20
1	AA	485	A	C5-C6-N6	-5.46	119.33	123.70
53	B5	1808	G	C5-C6-O6	-5.46	125.32	128.60
53	B5	2590	A	O4'-C1'-N9	5.46	112.57	108.20
53	B5	3150	A	O4'-C1'-N9	5.46	112.57	108.20
53	B5	3065	G	C5-C6-O6	-5.46	125.33	128.60
1	AA	551	G	O4'-C1'-N9	5.46	112.56	108.20
1	AA	824	G	C5-C6-O6	-5.46	125.33	128.60
1	AA	985	G	O4'-C1'-N9	5.46	112.56	108.20
1	AA	1240	A	C4-C5-C6	5.46	119.73	117.00
1	AA	1505	G	O4'-C1'-N9	5.46	112.56	108.20
52	B4	134	G	C5-C6-O6	-5.46	125.33	128.60
53	B5	550	A	O4'-C1'-N9	5.46	112.56	108.20
53	B5	2530	G	C5-C6-O6	-5.46	125.33	128.60
53	B5	2850	G	C5-C6-O6	-5.46	125.33	128.60
1	AA	885	G	O4'-C1'-N9	5.46	112.56	108.20
53	B5	347	G	O4'-C1'-N9	5.46	112.56	108.20
19	A7	15	G	C5-N7-C8	5.45	107.03	104.30
19	A7	20	G	C5-N7-C8	-5.45	101.57	104.30
53	B5	56	G	C5-C6-O6	-5.45	125.33	128.60
53	B5	143	G	C5-C6-O6	-5.45	125.33	128.60
53	B5	934	G	C5-C6-O6	-5.45	125.33	128.60
53	B5	1908	A	O4'-C1'-N9	5.45	112.56	108.20
53	B5	2690	G	C5-C6-O6	-5.45	125.33	128.60
19	A7	76	A	C1'-O4'-C4'	-5.45	105.54	109.90
53	B5	3099	C	N3-C4-C5	-5.45	119.72	121.90
1	AA	461	G	O4'-C1'-N9	5.45	112.56	108.20
1	AA	1259	G	C5-C6-O6	-5.45	125.33	128.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	956	U	O4'-C1'-N1	5.45	112.56	108.20
53	B5	1431	G	C5-C6-O6	-5.45	125.33	128.60
53	B5	3144	G	O4'-C1'-N9	5.45	112.56	108.20
53	B5	3348	G	O4'-C1'-N9	5.45	112.56	108.20
1	AA	23	G	O4'-C1'-N9	5.45	112.56	108.20
1	AA	1429	U	O4'-C1'-N1	5.45	112.56	108.20
33	BI	159	PHE	CB-CG-CD1	5.45	124.61	120.80
53	B5	353	G	C5-C6-O6	-5.45	125.33	128.60
53	B5	433	A	C4-C5-C6	5.45	119.72	117.00
53	B5	711	A	C5-C6-N6	-5.45	119.34	123.70
53	B5	2201	G	O4'-C1'-N9	5.45	112.56	108.20
53	B5	2636	A	C5-C6-N6	-5.45	119.34	123.70
53	B5	2957	G	C5-C6-O6	-5.45	125.33	128.60
1	AA	1124	G	O4'-C1'-N9	5.45	112.56	108.20
53	B5	187	A	C4-C5-C6	5.45	119.72	117.00
53	B5	575	G	C5-C6-O6	-5.45	125.33	128.60
53	B5	1780	G	O4'-C1'-N9	5.45	112.56	108.20
53	B5	2129	U	O4'-C1'-N1	5.45	112.56	108.20
53	B5	2136	C	O4'-C1'-N1	5.45	112.56	108.20
1	AA	422	G	C5-C6-O6	-5.45	125.33	128.60
1	AA	655	G	N1-C6-O6	5.45	123.17	119.90
1	AA	1230	A	O4'-C1'-N9	5.45	112.56	108.20
1	AA	1683	G	O4'-C1'-N9	5.45	112.56	108.20
19	A7	72	C	O3'-P-O5'	5.45	114.35	104.00
53	B5	585	A	O4'-C1'-N9	5.45	112.56	108.20
53	B5	843	A	O4'-C1'-N9	5.45	112.56	108.20
1	AA	83	G	O4'-C1'-N9	5.44	112.56	108.20
1	AA	127	G	C5-C6-O6	-5.44	125.33	128.60
1	AA	171	A	O4'-C1'-N9	5.44	112.56	108.20
53	B5	1266	G	C5-C6-O6	-5.44	125.33	128.60
1	AA	100	A	C4-C5-C6	5.44	119.72	117.00
1	AA	1473	A	C5-C6-N6	-5.44	119.35	123.70
27	BC	116	ARG	NE-CZ-NH1	5.44	123.02	120.30
53	B5	2608	G	N1-C6-O6	5.44	123.17	119.90
53	B5	3232	G	C5-C6-O6	-5.44	125.33	128.60
1	AA	465	G	O4'-C1'-N9	5.44	112.55	108.20
53	B5	394	G	O4'-C1'-N9	5.44	112.55	108.20
1	AA	278	U	O4'-C1'-N1	5.44	112.55	108.20
1	AA	624	G	C5-C6-O6	-5.44	125.34	128.60
51	B3	96	G	O4'-C1'-N9	5.44	112.55	108.20
53	B5	3337	G	O4'-C1'-N9	5.44	112.55	108.20
53	B5	1394	A	O4'-C1'-N9	5.44	112.55	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	1715	A	O4'-C1'-N9	5.44	112.55	108.20
53	B5	2592	G	O4'-C1'-N9	5.44	112.55	108.20
1	AA	1212	C	O4'-C1'-N1	5.44	112.55	108.20
1	AA	1338	A	O4'-C1'-N9	5.44	112.55	108.20
35	BK	121	PHE	CB-CG-CD2	-5.44	117.00	120.80
53	B5	72	C	N3-C4-N4	5.44	121.81	118.00
53	B5	692	A	O4'-C1'-N9	5.44	112.55	108.20
53	B5	2279	A	O4'-C1'-N9	5.44	112.55	108.20
53	B5	3310	A	C4-C5-C6	5.44	119.72	117.00
1	AA	373	G	O4'-C1'-N9	5.43	112.55	108.20
53	B5	331	G	O4'-C1'-N9	5.43	112.55	108.20
53	B5	374	A	O4'-C1'-N9	5.43	112.55	108.20
53	B5	932	U	P-O3'-C3'	5.43	126.22	119.70
53	B5	1048	A	C4-C5-C6	5.43	119.72	117.00
53	B5	1097	G	C5-C6-O6	-5.43	125.34	128.60
53	B5	1915	A	C5-C6-N6	-5.43	119.35	123.70
53	B5	3268	G	C5-C6-O6	-5.43	125.34	128.60
1	AA	283	U	O4'-C1'-N1	5.43	112.55	108.20
1	AA	1133	U	O4'-C1'-N1	5.43	112.55	108.20
1	AA	1135	A	O4'-C1'-N9	5.43	112.55	108.20
1	AA	1752	A	O4'-C1'-N9	5.43	112.55	108.20
19	A7	68	U	C4-C5-C6	5.43	122.96	119.70
53	B5	239	G	C5-C6-O6	-5.43	125.34	128.60
53	B5	1066	G	O4'-C1'-N9	5.43	112.55	108.20
53	B5	2902	A	C5-C6-N6	-5.43	119.35	123.70
53	B5	3054	U	O4'-C1'-N1	5.43	112.55	108.20
53	B5	2145	A	C4-C5-C6	5.43	119.72	117.00
1	AA	17	C	N3-C4-N4	5.43	121.80	118.00
1	AA	62	A	C5-C6-N6	-5.43	119.36	123.70
1	AA	1040	A	P-O3'-C3'	5.43	126.21	119.70
1	AA	1049	U	P-O3'-C3'	5.43	126.22	119.70
1	AA	1061	G	N1-C6-O6	5.43	123.16	119.90
1	AA	1660	G	O4'-C1'-N9	5.43	112.54	108.20
53	B5	350	C	N3-C4-C5	-5.43	119.73	121.90
53	B5	740	G	O4'-C1'-N9	5.43	112.54	108.20
53	B5	1113	G	C5-C6-O6	-5.43	125.34	128.60
53	B5	2192	C	N3-C4-N4	5.43	121.80	118.00
53	B5	2840	C	O4'-C1'-N1	5.43	112.54	108.20
53	B5	3188	G	O4'-C1'-N9	5.43	112.54	108.20
53	B5	3245	A	C4-C5-C6	5.43	119.72	117.00
1	AA	1646	A	O4'-C1'-N9	5.43	112.54	108.20
19	A7	73	A	C5-C6-N1	5.43	120.41	117.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	117	U	O4'-C1'-N1	5.43	112.54	108.20
1	AA	1653	A	O4'-C1'-N9	5.43	112.54	108.20
19	A7	6	U	C4'-C3'-C2'	-5.43	97.17	102.60
19	A7	45	G	C5'-C4'-C3'	-5.43	107.32	116.00
53	B5	538	G	O4'-C1'-N9	5.43	112.54	108.20
53	B5	617	G	O4'-C1'-N9	5.43	112.54	108.20
53	B5	1519	G	O4'-C1'-N9	5.43	112.54	108.20
53	B5	3142	A	O4'-C1'-N9	5.43	112.54	108.20
1	AA	262	U	O4'-C1'-N1	5.42	112.54	108.20
1	AA	608	U	P-O3'-C3'	5.42	126.21	119.70
1	AA	1754	A	O4'-C1'-N9	5.42	112.54	108.20
1	AA	1784	G	O4'-C1'-N9	5.42	112.54	108.20
29	BE	187	THR	C-N-CA	5.42	135.26	121.70
53	B5	1536	G	C5-C6-O6	-5.42	125.34	128.60
53	B5	3128	G	O4'-C1'-N9	5.42	112.54	108.20
53	B5	103	G	O4'-C1'-N9	5.42	112.54	108.20
53	B5	3294	A	O4'-C1'-N9	5.42	112.54	108.20
1	AA	47	A	C4-C5-C6	5.42	119.71	117.00
1	AA	953	G	C5-C6-O6	-5.42	125.35	128.60
1	AA	1035	A	O4'-C1'-N9	5.42	112.54	108.20
1	AA	1522	A	C4'-C3'-C2'	5.42	108.02	102.60
1	AA	1522	A	N1-C6-N6	5.42	121.85	118.60
1	AA	1786	G	O4'-C1'-N9	5.42	112.54	108.20
19	A7	48	C	N3-C2-O2	-5.42	118.11	121.90
53	B5	719	U	O4'-C1'-N1	5.42	112.54	108.20
53	B5	3086	A	C5-C6-N6	-5.42	119.36	123.70
1	AA	913	G	O4'-C1'-N9	5.42	112.54	108.20
1	AA	1760	A	C5-C6-N6	-5.42	119.36	123.70
53	B5	2937	G	C5-C6-O6	-5.42	125.35	128.60
1	AA	213	A	O4'-C1'-N9	5.42	112.53	108.20
1	AA	1226	A	C5-C6-N6	-5.42	119.36	123.70
1	AA	1239	G	C5-C6-O6	-5.42	125.35	128.60
1	AA	1357	A	O4'-C1'-N9	5.42	112.53	108.20
9	AJ	72	ASN	N-CA-CB	5.42	120.35	110.60
53	B5	699	A	O4'-C1'-N9	5.42	112.53	108.20
53	B5	888	A	O4'-C1'-N9	5.42	112.53	108.20
53	B5	1093	A	O4'-C1'-N9	5.42	112.53	108.20
53	B5	1848	G	C5-C6-O6	-5.42	125.35	128.60
53	B5	2230	C	O4'-C1'-N1	5.42	112.53	108.20
53	B5	2337	C	N3-C4-N4	5.42	121.79	118.00
53	B5	2796	G	C5-C6-O6	-5.42	125.35	128.60
1	AA	196	G	O4'-C1'-N9	5.42	112.53	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1426	G	C5-C6-O6	-5.42	125.35	128.60
1	AA	1510	G	C5-C6-O6	-5.42	125.35	128.60
53	B5	1157	G	O4'-C1'-N9	5.42	112.53	108.20
1	AA	1116	G	O4'-C1'-N9	5.42	112.53	108.20
52	B4	72	G	C5-C6-O6	-5.42	125.35	128.60
53	B5	709	A	O4'-C1'-N9	5.42	112.53	108.20
53	B5	1862	U	O4'-C1'-N1	5.42	112.53	108.20
1	AA	78	A	O4'-C1'-N9	5.41	112.53	108.20
1	AA	397	A	O4'-C1'-N9	5.41	112.53	108.20
1	AA	1491	A	C4-C5-C6	5.41	119.71	117.00
52	B4	112	U	C2-N1-C1'	5.41	124.20	117.70
53	B5	58	G	C5-C6-O6	-5.41	125.35	128.60
53	B5	1829	G	O4'-C1'-N9	5.41	112.53	108.20
53	B5	3340	G	O4'-C1'-N9	5.41	112.53	108.20
1	AA	389	G	O4'-C1'-N9	5.41	112.53	108.20
53	B5	775	A	O4'-C1'-N9	5.41	112.53	108.20
53	B5	2607	G	C5-C6-O6	-5.41	125.35	128.60
1	AA	865	A	N9-C4-C5	5.41	107.96	105.80
1	AA	1614	G	C5-C6-O6	-5.41	125.35	128.60
4	AD	79	ARG	NE-CZ-NH1	5.41	123.00	120.30
53	B5	1017	C	N3-C4-N4	5.41	121.79	118.00
53	B5	2528	G	O4'-C1'-N9	5.41	112.53	108.20
1	AA	628	G	C5-C6-O6	-5.41	125.36	128.60
19	A7	69	U	O5'-C5'-C4'	5.41	121.98	111.70
53	B5	481	U	O4'-C1'-N1	5.41	112.53	108.20
53	B5	557	A	O4'-C1'-N9	5.41	112.53	108.20
53	B5	754	G	C5-C6-O6	-5.41	125.36	128.60
53	B5	1813	A	C4-C5-C6	5.41	119.70	117.00
53	B5	2188	A	O4'-C1'-N9	5.41	112.53	108.20
53	B5	2412	G	O4'-C1'-N9	5.41	112.53	108.20
1	AA	1090	A	O4'-C1'-N9	5.41	112.53	108.20
53	B5	808	A	C5-C6-N6	-5.41	119.37	123.70
53	B5	1113	G	O4'-C1'-N9	5.41	112.53	108.20
1	AA	326	G	O4'-C1'-N9	5.41	112.52	108.20
1	AA	624	G	O4'-C1'-N9	5.41	112.52	108.20
1	AA	952	G	O4'-C1'-N9	5.41	112.53	108.20
52	B4	152	G	C5-C6-O6	-5.41	125.36	128.60
53	B5	733	G	O4'-C1'-N9	5.41	112.53	108.20
53	B5	869	G	C5-C6-O6	-5.41	125.36	128.60
53	B5	2362	C	N3-C4-N4	5.41	121.78	118.00
1	AA	457	G	C5-C6-O6	-5.40	125.36	128.60
53	B5	901	G	N1-C6-O6	5.40	123.14	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	1860	G	O4'-C1'-N9	5.40	112.52	108.20
1	AA	366	A	O4'-C1'-N9	5.40	112.52	108.20
1	AA	901	G	C5-C6-O6	-5.40	125.36	128.60
1	AA	1280	U	O4'-C1'-N1	5.40	112.52	108.20
53	B5	308	A	O4'-C1'-N9	5.40	112.52	108.20
53	B5	579	G	C5-C6-O6	-5.40	125.36	128.60
53	B5	2316	G	O4'-C1'-N9	5.40	112.52	108.20
53	B5	3096	C	N3-C4-C5	-5.40	119.74	121.90
1	AA	1742	A	O4'-C1'-N9	5.40	112.52	108.20
53	B5	258	G	O4'-C1'-N9	5.40	112.52	108.20
19	A7	63	C	C5-C6-N1	5.40	123.70	121.00
52	B4	136	G	C5-C6-O6	-5.40	125.36	128.60
53	B5	141	C	P-O3'-C3'	5.40	126.18	119.70
53	B5	630	A	O4'-C1'-N9	5.40	112.52	108.20
53	B5	3292	A	C4-C5-C6	5.40	119.70	117.00
53	B5	864	G	O4'-C1'-N9	5.40	112.52	108.20
53	B5	1372	C	N3-C4-N4	5.40	121.78	118.00
53	B5	2425	G	O4'-C1'-N9	5.40	112.52	108.20
1	AA	865	A	C8-N9-C4	-5.39	103.64	105.80
53	B5	22	G	O4'-C1'-N9	5.39	112.52	108.20
53	B5	1025	A	O4'-C1'-N9	5.39	112.52	108.20
53	B5	1307	G	C5-C6-O6	-5.39	125.36	128.60
53	B5	2171	G	C5-C6-O6	-5.39	125.36	128.60
53	B5	2224	A	C4-C5-C6	5.39	119.70	117.00
1	AA	802	G	N1-C6-O6	5.39	123.14	119.90
1	AA	1209	G	N1-C6-O6	5.39	123.14	119.90
53	B5	932	U	O4'-C1'-N1	5.39	112.51	108.20
53	B5	1281	G	C5-C6-O6	-5.39	125.36	128.60
53	B5	2387	A	C4-C5-C6	5.39	119.70	117.00
1	AA	325	G	N1-C6-O6	5.39	123.14	119.90
1	AA	604	A	O4'-C1'-N9	5.39	112.51	108.20
1	AA	1123	G	O4'-C1'-N9	5.39	112.51	108.20
53	B5	2563	G	O4'-C1'-N9	5.39	112.51	108.20
53	B5	2745	G	O4'-C1'-N9	5.39	112.51	108.20
1	AA	953	G	O4'-C1'-N9	5.39	112.51	108.20
1	AA	1185	A	C5-C6-N6	-5.39	119.39	123.70
1	AA	1690	G	O4'-C1'-N9	5.39	112.51	108.20
52	B4	31	G	C5-C6-O6	-5.39	125.37	128.60
53	B5	937	G	C5-C6-O6	-5.39	125.37	128.60
1	AA	62	A	C4-C5-C6	5.39	119.69	117.00
1	AA	462	G	O4'-C1'-N9	5.39	112.51	108.20
1	AA	526	A	C4-C5-C6	5.39	119.69	117.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	23	A	O4'-C1'-N9	5.39	112.51	108.20
53	B5	385	A	C4-C5-C6	5.39	119.69	117.00
53	B5	3191	G	C5-C6-O6	-5.39	125.37	128.60
1	AA	756	A	C5-C6-N6	-5.39	119.39	123.70
1	AA	911	U	O4'-C1'-N1	5.39	112.51	108.20
1	AA	1083	A	O4'-C1'-N9	5.39	112.51	108.20
1	AA	1232	A	O4'-C1'-N9	5.39	112.51	108.20
52	B4	14	C	N3-C4-N4	5.39	121.77	118.00
53	B5	357	A	O4'-C1'-N9	5.39	112.51	108.20
53	B5	836	A	C5-C6-N6	-5.39	119.39	123.70
53	B5	1477	A	O4'-C1'-N9	5.39	112.51	108.20
53	B5	2215	A	C4-C5-C6	5.39	119.69	117.00
1	AA	151	G	O4'-C1'-N9	5.38	112.51	108.20
1	AA	878	G	N1-C6-O6	5.38	123.13	119.90
1	AA	1316	A	O4'-C1'-N9	5.38	112.51	108.20
53	B5	42	C	N3-C4-N4	5.38	121.77	118.00
53	B5	1026	A	O4'-C1'-N9	5.38	112.51	108.20
53	B5	2195	C	N3-C4-C5	-5.38	119.75	121.90
53	B5	2430	A	O4'-C1'-N9	5.38	112.51	108.20
1	AA	561	G	O4'-C1'-N9	5.38	112.50	108.20
1	AA	876	G	O4'-C1'-N9	5.38	112.50	108.20
1	AA	975	G	C5-C6-O6	-5.38	125.37	128.60
1	AA	1376	C	O4'-C1'-N1	5.38	112.51	108.20
19	A7	1	G	N1-C2-N3	5.38	127.13	123.90
19	A7	11	C	N1-C2-N3	5.38	122.97	119.20
53	B5	1769	G	C5-C6-O6	-5.38	125.37	128.60
1	AA	1128	A	O4'-C1'-N9	5.38	112.50	108.20
1	AA	1184	G	O4'-C1'-N9	5.38	112.50	108.20
53	B5	936	A	P-O3'-C3'	5.38	126.16	119.70
53	B5	2139	A	C5-C6-N6	-5.38	119.40	123.70
53	B5	2837	A	O4'-C1'-N9	5.38	112.50	108.20
53	B5	3362	A	C5-C6-N6	-5.38	119.40	123.70
1	AA	1233	G	O4'-C1'-N9	5.38	112.50	108.20
1	AA	1541	A	C4-C5-C6	5.38	119.69	117.00
53	B5	1492	G	O4'-C1'-N9	5.38	112.50	108.20
53	B5	2773	C	C2-N1-C1'	5.38	124.72	118.80
19	A7	5	A	C4-C5-N7	5.38	113.39	110.70
53	B5	1027	A	C5-C6-N6	-5.38	119.40	123.70
53	B5	1162	U	P-O5'-C5'	-5.38	112.30	120.90
53	B5	2312	A	C5-C6-N6	-5.38	119.40	123.70
1	AA	987	A	O4'-C1'-N9	5.38	112.50	108.20
53	B5	319	A	C4-C5-C6	5.38	119.69	117.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	373	A	C5-C6-N1	-5.38	115.01	117.70
53	B5	390	G	P-O3'-C3'	5.38	126.15	119.70
53	B5	773	G	O4'-C1'-N9	5.38	112.50	108.20
53	B5	1367	G	C5-C6-O6	-5.38	125.38	128.60
1	AA	1089	A	C4-C5-C6	5.37	119.69	117.00
1	AA	1505	G	C5-C6-O6	-5.37	125.38	128.60
19	A7	5	A	C2-N3-C4	5.37	113.29	110.60
51	B3	37	G	C5-C6-O6	-5.37	125.38	128.60
52	B4	37	A	C4-C5-C6	5.37	119.69	117.00
53	B5	474	G	C5-C6-O6	-5.37	125.38	128.60
53	B5	1884	A	O4'-C1'-N9	5.37	112.50	108.20
53	B5	2956	A	C5-C6-N6	-5.37	119.40	123.70
1	AA	924	G	O4'-C1'-N9	5.37	112.50	108.20
1	AA	938	A	C5-C6-N6	-5.37	119.40	123.70
1	AA	1515	U	O4'-C1'-N1	5.37	112.50	108.20
1	AA	1518	U	O4'-C1'-N1	5.37	112.50	108.20
53	B5	862	U	O4'-C1'-N1	5.37	112.50	108.20
53	B5	1817	G	C5-C6-O6	-5.37	125.38	128.60
53	B5	2149	A	C5-C6-N6	-5.37	119.40	123.70
53	B5	2172	A	C4-C5-C6	5.37	119.69	117.00
53	B5	2488	A	O4'-C1'-N9	5.37	112.50	108.20
53	B5	2903	A	O4'-C1'-N9	5.37	112.50	108.20
53	B5	3080	G	O4'-C1'-N9	5.37	112.50	108.20
53	B5	1914	G	O4'-C1'-N9	5.37	112.50	108.20
1	AA	219	A	O4'-C1'-N9	5.37	112.50	108.20
1	AA	951	A	C5-C6-N6	-5.37	119.41	123.70
1	AA	1115	G	C5-C6-O6	-5.37	125.38	128.60
1	AA	1257	G	O4'-C1'-N9	5.37	112.49	108.20
1	AA	1501	A	O4'-C1'-N9	5.37	112.50	108.20
51	B3	89	A	O4'-C1'-N9	5.37	112.50	108.20
53	B5	938	C	N3-C4-N4	5.37	121.76	118.00
53	B5	2260	U	O4'-C1'-N1	5.37	112.50	108.20
53	B5	2398	A	O4'-C1'-N9	5.37	112.50	108.20
53	B5	761	A	O4'-C1'-N9	5.37	112.49	108.20
53	B5	1213	G	C5-C6-O6	-5.37	125.38	128.60
53	B5	2185	G	O4'-C1'-N9	5.37	112.49	108.20
1	AA	885	G	C5-C6-O6	-5.37	125.38	128.60
1	AA	1007	G	C5-C6-O6	-5.37	125.38	128.60
1	AA	1060	U	O4'-C1'-N1	5.37	112.49	108.20
1	AA	1573	G	O4'-C1'-N9	5.37	112.49	108.20
1	AA	1700	A	C4-C5-C6	5.37	119.68	117.00
19	A7	2	C	C2-N3-C4	-5.37	117.22	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
19	A7	9	A	C6-C5-N7	5.37	136.06	132.30
53	B5	522	A	O4'-C1'-N9	5.37	112.49	108.20
53	B5	1503	A	O4'-C1'-N9	5.37	112.49	108.20
53	B5	3044	G	O4'-C1'-N9	5.37	112.49	108.20
1	AA	163	G	O4'-C1'-N9	5.36	112.49	108.20
1	AA	243	G	C5-C6-O6	-5.36	125.38	128.60
1	AA	445	A	C5-C6-N6	-5.36	119.41	123.70
1	AA	905	A	O4'-C1'-N9	5.36	112.49	108.20
1	AA	972	A	O4'-C1'-N9	5.36	112.49	108.20
1	AA	1395	U	O4'-C1'-N1	5.36	112.49	108.20
52	B4	49	G	O4'-C1'-N9	5.36	112.49	108.20
53	B5	358	G	C5-C6-O6	-5.36	125.38	128.60
53	B5	1461	A	O4'-C1'-N9	5.36	112.49	108.20
53	B5	2730	G	O4'-C1'-N9	5.36	112.49	108.20
53	B5	3172	G	O4'-C1'-N9	5.36	112.49	108.20
53	B5	3285	C	N3-C4-N4	5.36	121.75	118.00
1	AA	460	A	O4'-C1'-N9	5.36	112.49	108.20
1	AA	1208	G	C5-C6-O6	-5.36	125.38	128.60
1	AA	1478	G	C5-C6-O6	-5.36	125.38	128.60
19	A7	2	C	C5-C4-N4	5.36	123.95	120.20
53	B5	2106	A	O4'-C1'-N9	5.36	112.49	108.20
53	B5	2805	G	O4'-C1'-N9	5.36	112.49	108.20
51	B3	46	A	C5'-C4'-C3'	5.36	124.58	116.00
52	B4	109	A	O4'-C1'-N9	5.36	112.49	108.20
53	B5	184	U	O4'-C1'-N1	5.36	112.49	108.20
53	B5	900	G	C5-C6-O6	-5.36	125.38	128.60
1	AA	1327	G	C5-C6-O6	-5.36	125.39	128.60
1	AA	1522	A	C1'-O4'-C4'	-5.36	105.61	109.90
1	AA	41	A	O4'-C1'-N9	5.36	112.49	108.20
1	AA	702	G	C5-C6-O6	-5.36	125.39	128.60
29	BE	22	ARG	NE-CZ-NH1	5.36	122.98	120.30
52	B4	22	U	O4'-C1'-N1	5.36	112.49	108.20
53	B5	1462	A	O4'-C1'-N9	5.36	112.49	108.20
53	B5	1510	G	C5-C6-O6	-5.36	125.39	128.60
53	B5	1901	A	P-O3'-C3'	5.36	126.13	119.70
53	B5	3039	C	N3-C4-C5	-5.36	119.76	121.90
1	AA	315	A	O4'-C1'-N9	5.36	112.48	108.20
19	A7	74	C	C4-C5-C6	5.36	120.08	117.40
53	B5	1419	A	C5-C6-N6	-5.36	119.42	123.70
53	B5	2161	G	O4'-C1'-N9	5.36	112.48	108.20
53	B5	2601	A	O4'-C1'-N9	5.36	112.48	108.20
53	B5	2656	A	C4-C5-C6	5.36	119.68	117.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	2924	U	O4'-C1'-N1	5.36	112.48	108.20
1	AA	1222	A	O4'-C1'-N9	5.35	112.48	108.20
53	B5	5	G	O4'-C1'-N9	5.35	112.48	108.20
53	B5	1781	C	N3-C4-N4	5.35	121.75	118.00
53	B5	2331	C	N3-C4-C5	-5.35	119.76	121.90
1	AA	1319	A	C4-C5-C6	5.35	119.68	117.00
53	B5	602	A	C5-C6-N6	-5.35	119.42	123.70
53	B5	2234	G	O4'-C1'-N9	5.35	112.48	108.20
53	B5	2632	G	C5-C6-O6	-5.35	125.39	128.60
53	B5	3143	C	O4'-C1'-N1	5.35	112.48	108.20
1	AA	1730	A	C5-C6-N6	-5.35	119.42	123.70
53	B5	1622	U	O4'-C1'-N1	5.35	112.48	108.20
1	AA	279	G	C5-C6-O6	-5.35	125.39	128.60
1	AA	1192	A	O4'-C1'-N9	5.35	112.48	108.20
1	AA	1445	C	O4'-C1'-N1	5.35	112.48	108.20
1	AA	1550	U	C1'-C2'-O2'	5.35	126.64	110.60
1	AA	1761	A	O4'-C1'-N9	5.35	112.48	108.20
51	B3	8	G	O4'-C1'-N9	5.35	112.48	108.20
53	B5	254	A	C5-C6-N6	-5.35	119.42	123.70
53	B5	1661	G	C5-C6-O6	-5.35	125.39	128.60
53	B5	1835	A	C4-C5-C6	5.35	119.67	117.00
53	B5	3099	C	O4'-C1'-N1	5.35	112.48	108.20
53	B5	1898	G	C5-C6-O6	-5.35	125.39	128.60
53	B5	2897	A	C4-C5-C6	5.35	119.67	117.00
1	AA	96	G	O4'-C1'-N9	5.34	112.47	108.20
1	AA	687	G	O4'-C1'-N9	5.34	112.47	108.20
1	AA	940	A	C4-C5-C6	5.34	119.67	117.00
52	B4	44	A	C5-C6-N6	-5.34	119.42	123.70
53	B5	62	A	O4'-C1'-N9	5.34	112.48	108.20
53	B5	157	A	O4'-C1'-N9	5.34	112.48	108.20
53	B5	303	G	N1-C6-O6	5.34	123.11	119.90
53	B5	937	G	O4'-C1'-N9	5.34	112.48	108.20
53	B5	1203	A	C5-C6-N6	-5.34	119.42	123.70
53	B5	1306	G	O4'-C1'-N9	5.34	112.47	108.20
53	B5	1547	G	O4'-C1'-N9	5.34	112.47	108.20
53	B5	1634	G	C5-C6-O6	-5.34	125.39	128.60
53	B5	2251	G	N1-C6-O6	5.34	123.11	119.90
53	B5	3367	C	N3-C4-C5	-5.34	119.76	121.90
1	AA	367	A	O4'-C1'-N9	5.34	112.47	108.20
1	AA	1700	A	O4'-C1'-N9	5.34	112.47	108.20
53	B5	2645	G	C5-C6-O6	-5.34	125.39	128.60
53	B5	2943	G	O4'-C1'-N9	5.34	112.47	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	3065	G	O4'-C1'-N9	5.34	112.47	108.20
1	AA	459	G	C5-C6-O6	-5.34	125.40	128.60
53	B5	373	A	O4'-C1'-N9	5.34	112.47	108.20
53	B5	1260	A	O4'-C1'-N9	5.34	112.47	108.20
1	AA	1435	G	O4'-C1'-N9	5.34	112.47	108.20
53	B5	597	G	O4'-C1'-N9	5.34	112.47	108.20
53	B5	600	G	C5-C6-O6	-5.34	125.40	128.60
53	B5	1264	G	O4'-C1'-N9	5.34	112.47	108.20
53	B5	2486	A	C4-C5-C6	5.34	119.67	117.00
53	B5	2691	A	C5-C6-N6	-5.34	119.43	123.70
53	B5	1926	C	O4'-C1'-N1	5.34	112.47	108.20
53	B5	2830	G	C5'-C4'-O4'	5.34	115.50	109.10
1	AA	865	A	C1'-O4'-C4'	-5.34	105.63	109.90
53	B5	98	G	O4'-C1'-N9	5.34	112.47	108.20
53	B5	833	G	C5-C6-O6	-5.34	125.40	128.60
53	B5	1153	A	C4-C5-C6	5.34	119.67	117.00
53	B5	2265	C	N3-C4-C5	-5.34	119.77	121.90
53	B5	3127	A	O4'-C1'-N9	5.34	112.47	108.20
53	B5	1242	G	C5-C6-O6	-5.33	125.40	128.60
1	AA	1085	A	C5-C6-N6	-5.33	119.43	123.70
1	AA	1139	A	O4'-C1'-N9	5.33	112.47	108.20
19	A7	64	A	C5-C6-N6	5.33	127.97	123.70
51	B3	20	A	O4'-C1'-N9	5.33	112.47	108.20
53	B5	116	A	P-O3'-C3'	5.33	126.10	119.70
53	B5	844	G	O4'-C1'-N9	5.33	112.47	108.20
53	B5	1540	U	O4'-C1'-N1	5.33	112.47	108.20
51	B3	52	U	O4'-C1'-N1	5.33	112.46	108.20
53	B5	708	G	C5-C6-O6	-5.33	125.40	128.60
53	B5	1333	C	N3-C4-N4	5.33	121.73	118.00
53	B5	1365	G	O4'-C1'-N9	5.33	112.47	108.20
53	B5	2399	A	O4'-C1'-N9	5.33	112.47	108.20
53	B5	2720	G	O4'-C1'-N9	5.33	112.47	108.20
1	AA	173	A	P-O3'-C3'	5.33	126.10	119.70
1	AA	1078	A	O4'-C1'-N9	5.33	112.46	108.20
1	AA	1274	G	O4'-C1'-N9	5.33	112.46	108.20
51	B3	7	G	O4'-C1'-N9	5.33	112.46	108.20
53	B5	568	G	O4'-C1'-N9	5.33	112.46	108.20
53	B5	2705	A	O4'-C1'-N9	5.33	112.46	108.20
51	B3	59	G	C5-C6-O6	-5.33	125.40	128.60
53	B5	4	U	O4'-C1'-N1	5.33	112.46	108.20
53	B5	499	G	O4'-C1'-N9	5.33	112.46	108.20
53	B5	813	G	O4'-C1'-N9	5.33	112.46	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	1193	A	C5-C6-N6	-5.33	119.44	123.70
53	B5	1239	C	N3-C4-C5	-5.33	119.77	121.90
53	B5	1287	A	C5-C6-N6	-5.33	119.44	123.70
1	AA	1022	A	C4-C5-C6	5.33	119.66	117.00
53	B5	106	A	O4'-C1'-N9	5.33	112.46	108.20
53	B5	1804	A	P-O3'-C3'	5.33	126.09	119.70
1	AA	274	G	O4'-C1'-N9	5.33	112.46	108.20
1	AA	1361	G	O4'-C1'-N9	5.33	112.46	108.20
19	A7	70	C	O5'-P-OP2	-5.33	100.91	105.70
52	B4	114	G	P-O3'-C3'	5.33	126.09	119.70
53	B5	37	U	O4'-C1'-N1	5.33	112.46	108.20
53	B5	1025	A	C5-C6-N6	-5.33	119.44	123.70
53	B5	1101	G	C5-C6-O6	-5.33	125.41	128.60
53	B5	1516	C	N3-C4-N4	5.33	121.73	118.00
53	B5	2183	A	O4'-C1'-N9	5.33	112.46	108.20
53	B5	2222	A	O4'-C1'-N9	5.33	112.46	108.20
53	B5	2838	A	C5-C6-N6	-5.33	119.44	123.70
53	B5	3047	U	O4'-C1'-N1	5.33	112.46	108.20
1	AA	859	A	N9-C1'-C2'	5.32	120.92	114.00
53	B5	590	G	C5-C6-O6	-5.32	125.41	128.60
53	B5	1441	G	O4'-C1'-N9	5.32	112.46	108.20
53	B5	1504	A	O4'-C1'-N9	5.32	112.46	108.20
53	B5	2121	G	C5-C6-O6	-5.32	125.41	128.60
53	B5	2288	G	O4'-C1'-N9	5.32	112.46	108.20
1	AA	1136	A	O4'-C1'-N9	5.32	112.46	108.20
1	AA	1417	C	O4'-C1'-N1	5.32	112.46	108.20
53	B5	145	G	O4'-C1'-N9	5.32	112.46	108.20
53	B5	1440	G	C5-C6-O6	-5.32	125.41	128.60
1	AA	1383	G	C5-C6-O6	-5.32	125.41	128.60
52	B4	53	A	O4'-C1'-N9	5.32	112.46	108.20
53	B5	704	U	O4'-C1'-N1	5.32	112.46	108.20
53	B5	2528	G	C5-C6-O6	-5.32	125.41	128.60
53	B5	2592	G	C5-C6-O6	-5.32	125.41	128.60
53	B5	2977	G	O4'-C1'-N9	5.32	112.46	108.20
53	B5	1937	U	O4'-C1'-N1	5.32	112.45	108.20
1	AA	124	A	O4'-C1'-N9	5.32	112.45	108.20
53	B5	775	A	C5-C6-N6	-5.32	119.45	123.70
53	B5	830	A	O4'-C1'-N9	5.32	112.45	108.20
53	B5	1063	G	C5-C6-O6	-5.32	125.41	128.60
53	B5	1446	A	C5-C6-N6	-5.32	119.44	123.70
53	B5	1643	A	O4'-C1'-N9	5.32	112.45	108.20
53	B5	1910	A	O4'-C1'-N9	5.32	112.45	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	2439	A	O4'-C1'-N9	5.32	112.45	108.20
1	AA	1129	A	C4-C5-C6	5.32	119.66	117.00
1	AA	1327	G	O4'-C1'-N9	5.32	112.45	108.20
1	AA	16	G	O4'-C1'-N9	5.31	112.45	108.20
53	B5	303	G	O4'-C1'-N9	5.31	112.45	108.20
53	B5	3031	G	O4'-C1'-N9	5.31	112.45	108.20
1	AA	61	A	C5-C6-N6	-5.31	119.45	123.70
1	AA	71	A	O4'-C1'-N9	5.31	112.45	108.20
1	AA	72	A	O4'-C1'-N9	5.31	112.45	108.20
1	AA	671	G	O4'-C1'-N9	5.31	112.45	108.20
1	AA	941	G	C5-C6-O6	-5.31	125.41	128.60
1	AA	1176	G	O4'-C1'-N9	5.31	112.45	108.20
53	B5	52	A	C4-C5-C6	5.31	119.66	117.00
53	B5	160	G	O4'-C1'-N9	5.31	112.45	108.20
53	B5	201	A	C5-C6-N6	-5.31	119.45	123.70
53	B5	613	G	O4'-C1'-N9	5.31	112.45	108.20
53	B5	1604	G	C5-C6-O6	-5.31	125.41	128.60
53	B5	2198	A	O4'-C1'-N9	5.31	112.45	108.20
53	B5	3238	G	O4'-C1'-N9	5.31	112.45	108.20
1	AA	63	G	O4'-C1'-N9	5.31	112.45	108.20
1	AA	384	G	C5-C6-O6	-5.31	125.41	128.60
1	AA	776	G	O4'-C1'-N9	5.31	112.45	108.20
53	B5	2321	A	O4'-C1'-N9	5.31	112.45	108.20
53	B5	3361	G	O4'-C1'-N9	5.31	112.45	108.20
1	AA	615	A	C5-C6-N6	-5.31	119.45	123.70
1	AA	1548	A	C4-C5-C6	5.31	119.66	117.00
1	AA	1640	G	C5-C6-O6	-5.31	125.41	128.60
1	AA	1718	G	O4'-C1'-N9	5.31	112.45	108.20
1	AA	1755	G	C5-C6-O6	-5.31	125.42	128.60
53	B5	319	A	C5-C6-N1	-5.31	115.05	117.70
53	B5	2450	G	C5-C6-O6	-5.31	125.41	128.60
53	B5	2847	A	C4-C5-C6	5.31	119.66	117.00
53	B5	2848	G	P-O3'-C3'	5.31	126.07	119.70
53	B5	2848	G	N1-C6-O6	5.31	123.09	119.90
53	B5	3144	G	C5-C6-O6	-5.31	125.41	128.60
1	AA	631	G	O4'-C1'-N9	5.31	112.45	108.20
1	AA	900	G	O4'-C1'-N9	5.31	112.45	108.20
1	AA	1291	G	C5-C6-O6	-5.31	125.42	128.60
1	AA	1428	C	O4'-C1'-N1	5.31	112.45	108.20
19	A7	7	U	N1-C2-N3	5.31	118.08	114.90
53	B5	1079	A	C5-C6-N6	-5.31	119.45	123.70
53	B5	2791	G	C5-C6-O6	-5.31	125.42	128.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	3328	G	C5-C6-O6	-5.31	125.42	128.60
53	B5	319	A	O4'-C1'-N9	5.31	112.44	108.20
53	B5	2878	G	C5-C6-O6	-5.31	125.42	128.60
1	AA	81	G	O4'-C1'-N9	5.30	112.44	108.20
1	AA	1144	A	C4-C5-C6	5.30	119.65	117.00
51	B3	115	A	O4'-C1'-N9	5.30	112.44	108.20
53	B5	648	C	O4'-C1'-N1	5.30	112.44	108.20
53	B5	1136	A	O4'-C1'-N9	5.30	112.44	108.20
53	B5	1300	G	O4'-C1'-N9	5.30	112.44	108.20
53	B5	1385	C	N3-C4-N4	5.30	121.71	118.00
53	B5	1650	G	C5-C6-O6	-5.30	125.42	128.60
53	B5	3030	G	O4'-C1'-N9	5.30	112.44	108.20
1	AA	26	A	O4'-C1'-N9	5.30	112.44	108.20
1	AA	776	G	C5-C6-O6	-5.30	125.42	128.60
1	AA	92	A	C4-C5-C6	5.30	119.65	117.00
1	AA	809	A	O4'-C1'-N9	5.30	112.44	108.20
1	AA	1194	G	C5-C6-O6	-5.30	125.42	128.60
53	B5	707	U	O4'-C1'-N1	5.30	112.44	108.20
53	B5	997	A	C4-C5-C6	5.30	119.65	117.00
53	B5	2134	G	C5-C6-O6	-5.30	125.42	128.60
19	A7	9	A	N9-C1'-C2'	-5.30	106.17	112.00
53	B5	114	A	O4'-C1'-N9	5.30	112.44	108.20
53	B5	1344	G	C5-C6-O6	-5.30	125.42	128.60
53	B5	1380	G	N1-C6-O6	5.30	123.08	119.90
1	AA	488	G	C5-C6-O6	-5.30	125.42	128.60
53	B5	27	C	N3-C4-N4	5.30	121.71	118.00
1	AA	1421	A	O4'-C1'-N9	5.30	112.44	108.20
1	AA	1477	A	O4'-C1'-N9	5.30	112.44	108.20
29	BE	207	TYR	CB-CG-CD1	5.30	124.18	121.00
53	B5	196	G	O4'-C1'-N9	5.30	112.44	108.20
53	B5	234	G	C5-C6-O6	-5.30	125.42	128.60
53	B5	1165	A	C5-C6-N6	-5.30	119.46	123.70
53	B5	2575	G	N1-C6-O6	5.30	123.08	119.90
1	AA	1746	G	C5-C6-O6	-5.29	125.42	128.60
53	B5	1939	G	C5-C6-O6	-5.29	125.42	128.60
53	B5	2610	G	C5-C6-O6	-5.29	125.42	128.60
1	AA	215	A	O4'-C1'-N9	5.29	112.44	108.20
53	B5	476	G	O4'-C1'-N9	5.29	112.44	108.20
53	B5	700	C	N3-C4-N4	5.29	121.70	118.00
53	B5	831	G	C5-C6-O6	-5.29	125.42	128.60
53	B5	2366	C	N3-C4-N4	5.29	121.71	118.00
53	B5	3335	A	O4'-C1'-N9	5.29	112.44	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	470	A	O4'-C1'-N9	5.29	112.43	108.20
1	AA	684	A	O4'-C1'-N9	5.29	112.43	108.20
51	B3	54	A	C5-C6-N6	-5.29	119.47	123.70
53	B5	71	A	C4-C5-C6	5.29	119.65	117.00
53	B5	1155	C	N3-C4-N4	5.29	121.70	118.00
53	B5	2202	C	N3-C4-N4	5.29	121.70	118.00
53	B5	2901	G	O4'-C1'-N9	5.29	112.43	108.20
53	B5	3079	U	O4'-C1'-N1	5.29	112.43	108.20
1	AA	929	A	C5-C6-N6	-5.29	119.47	123.70
1	AA	935	G	C5-C6-O6	-5.29	125.43	128.60
1	AA	1726	A	O4'-C1'-N9	5.29	112.43	108.20
53	B5	1171	G	O4'-C1'-N9	5.29	112.43	108.20
53	B5	2437	G	O4'-C1'-N9	5.29	112.43	108.20
1	AA	214	G	C5-C6-O6	-5.29	125.43	128.60
1	AA	404	G	C5-C6-O6	-5.29	125.43	128.60
1	AA	610	G	C5-C6-O6	-5.29	125.43	128.60
52	B4	62	C	C6-N1-C1'	-5.29	114.45	120.80
53	B5	201	A	C4-C5-C6	5.29	119.64	117.00
53	B5	584	G	C5-C6-O6	-5.29	125.43	128.60
53	B5	1075	A	O4'-C1'-N9	5.29	112.43	108.20
53	B5	2687	G	O4'-C1'-N9	5.29	112.43	108.20
53	B5	2820	A	O4'-C1'-N9	5.29	112.43	108.20
53	B5	2911	A	O4'-C1'-N9	5.29	112.43	108.20
53	B5	3284	G	O4'-C1'-N9	5.29	112.43	108.20
1	AA	723	G	O4'-C1'-N9	5.29	112.43	108.20
53	B5	312	C	N3-C4-N4	5.29	121.70	118.00
53	B5	2454	G	C5-C6-O6	-5.29	125.43	128.60
1	AA	123	G	C5-C6-O6	-5.29	125.43	128.60
1	AA	669	G	O4'-C1'-N9	5.29	112.43	108.20
19	A7	21	A	C5'-C4'-C3'	-5.29	107.54	116.00
51	B3	24	A	O4'-C1'-N9	5.29	112.43	108.20
53	B5	1409	G	C5-C6-O6	-5.29	125.43	128.60
53	B5	1605	A	C5-C6-N6	-5.29	119.47	123.70
53	B5	1758	G	O4'-C1'-N9	5.29	112.43	108.20
53	B5	2612	U	O4'-C1'-N1	5.29	112.43	108.20
1	AA	647	G	C5-C6-O6	-5.28	125.43	128.60
1	AA	929	A	C4-C5-C6	5.28	119.64	117.00
1	AA	1087	C	N3-C4-N4	5.28	121.70	118.00
53	B5	15	C	N3-C4-N4	5.28	121.70	118.00
53	B5	109	A	O4'-C1'-N9	5.28	112.43	108.20
53	B5	264	G	O4'-C1'-N9	5.28	112.43	108.20
53	B5	274	G	O4'-C1'-N9	5.28	112.43	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	469	G	N1-C6-O6	5.28	123.07	119.90
53	B5	1182	A	C4-C5-C6	5.28	119.64	117.00
53	B5	1715	A	C5-C6-N1	-5.28	115.06	117.70
53	B5	1893	A	O4'-C1'-N9	5.28	112.43	108.20
53	B5	2515	A	C5-C6-N6	-5.28	119.47	123.70
53	B5	3027	A	O4'-C1'-N9	5.28	112.43	108.20
1	AA	1043	G	O4'-C1'-N9	5.28	112.43	108.20
1	AA	1605	G	C5-C6-O6	-5.28	125.43	128.60
53	B5	450	G	O4'-C1'-N9	5.28	112.42	108.20
53	B5	570	A	O4'-C1'-N9	5.28	112.42	108.20
53	B5	1079	A	C4-C5-C6	5.28	119.64	117.00
53	B5	1396	C	N3-C4-N4	5.28	121.70	118.00
53	B5	1618	G	C5-C6-O6	-5.28	125.43	128.60
1	AA	661	A	O4'-C1'-N9	5.28	112.42	108.20
19	A7	4	G	N3-C4-C5	5.28	131.24	128.60
51	B3	84	G	C5-C6-O6	-5.28	125.43	128.60
53	B5	5	G	C5-C6-O6	-5.28	125.43	128.60
53	B5	1879	A	C5-C6-N6	-5.28	119.48	123.70
1	AA	371	G	C5-C6-O6	-5.28	125.43	128.60
19	A7	25	C	O4'-C1'-N1	5.28	112.42	108.20
53	B5	3067	C	N3-C4-N4	5.28	121.69	118.00
1	AA	269	G	C5-C6-O6	-5.28	125.43	128.60
1	AA	384	G	O4'-C1'-N9	5.28	112.42	108.20
1	AA	399	A	O4'-C1'-N9	5.28	112.42	108.20
1	AA	1634	C	N1-C1'-C2'	5.28	120.86	114.00
53	B5	3101	G	O4'-C1'-N9	5.28	112.42	108.20
53	B5	3299	A	O4'-C1'-N9	5.28	112.42	108.20
53	B5	384	A	O4'-C1'-N9	5.28	112.42	108.20
53	B5	656	A	O4'-C1'-N9	5.28	112.42	108.20
53	B5	1932	A	C5-C6-N6	-5.28	119.48	123.70
53	B5	2208	A	O4'-C1'-N9	5.28	112.42	108.20
53	B5	2243	A	O4'-C1'-N9	5.28	112.42	108.20
53	B5	3279	A	C5-C6-N6	-5.28	119.48	123.70
1	AA	40	A	O4'-C1'-N9	5.27	112.42	108.20
1	AA	1007	G	O4'-C1'-N9	5.27	112.42	108.20
19	A7	18	G	N3-C4-C5	5.27	131.24	128.60
53	B5	703	G	C5-C6-O6	-5.27	125.44	128.60
53	B5	1760	A	O4'-C1'-N9	5.27	112.42	108.20
1	AA	651	G	O4'-C1'-N9	5.27	112.42	108.20
19	A7	44	A	C5-C6-N1	5.27	120.34	117.70
53	B5	17	G	O4'-C1'-N9	5.27	112.42	108.20
53	B5	89	A	C5-C6-N6	-5.27	119.48	123.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	1071	U	O4'-C1'-N1	5.27	112.42	108.20
53	B5	1165	A	C4-C5-C6	5.27	119.64	117.00
53	B5	1648	A	O4'-C1'-N9	5.27	112.42	108.20
53	B5	3122	A	C4-C5-C6	5.27	119.64	117.00
53	B5	3316	A	O4'-C1'-N9	5.27	112.42	108.20
1	AA	729	G	O4'-C1'-N9	5.27	112.42	108.20
53	B5	931	C	N3-C4-N4	5.27	121.69	118.00
53	B5	2384	A	O4'-C1'-N9	5.27	112.42	108.20
53	B5	2523	A	O4'-C1'-N9	5.27	112.42	108.20
53	B5	3239	G	O4'-C1'-N9	5.27	112.42	108.20
1	AA	251	A	O4'-C1'-N9	5.27	112.42	108.20
1	AA	754	A	O4'-C1'-N9	5.27	112.42	108.20
52	B4	150	G	O4'-C1'-N9	5.27	112.42	108.20
53	B5	1131	G	O4'-C1'-N9	5.27	112.42	108.20
53	B5	1291	A	O4'-C1'-N9	5.27	112.42	108.20
53	B5	2557	G	O4'-C1'-N9	5.27	112.42	108.20
53	B5	71	A	C5-C6-N1	-5.27	115.07	117.70
53	B5	594	A	O4'-C1'-N9	5.27	112.41	108.20
53	B5	692	A	C4-C5-C6	5.27	119.63	117.00
53	B5	1163	A	C5-C6-N6	-5.27	119.48	123.70
53	B5	2684	C	N3-C4-C5	-5.27	119.79	121.90
1	AA	1240	A	C1'-O4'-C4'	-5.27	105.69	109.90
1	AA	1263	G	C5-C6-O6	-5.27	125.44	128.60
1	AA	1519	G	O4'-C1'-N9	5.27	112.41	108.20
1	AA	1446	G	O4'-C1'-N9	5.26	112.41	108.20
1	AA	1458	A	O4'-C1'-N9	5.26	112.41	108.20
52	B4	44	A	C4-C5-C6	5.26	119.63	117.00
53	B5	31	C	N3-C4-C5	-5.26	119.79	121.90
53	B5	393	U	P-O3'-C3'	5.26	126.02	119.70
53	B5	1226	G	C5-C6-O6	-5.26	125.44	128.60
53	B5	1404	G	C5-C6-O6	-5.26	125.44	128.60
53	B5	2445	A	O4'-C1'-N9	5.26	112.41	108.20
53	B5	2816	G	C5-C6-O6	-5.26	125.44	128.60
1	AA	1365	G	O4'-C1'-N9	5.26	112.41	108.20
53	B5	909	G	C5-C6-O6	-5.26	125.44	128.60
1	AA	545	A	O4'-C1'-N9	5.26	112.41	108.20
1	AA	664	U	O4'-C1'-N1	5.26	112.41	108.20
53	B5	95	A	O4'-C1'-N9	5.26	112.41	108.20
53	B5	859	G	C5-C6-O6	-5.26	125.44	128.60
53	B5	1688	U	O4'-C1'-N1	5.26	112.41	108.20
53	B5	1745	C	P-O3'-C3'	5.26	126.01	119.70
53	B5	2239	G	O4'-C1'-N9	5.26	112.41	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1041	U	O4'-C1'-N1	5.26	112.41	108.20
53	B5	1461	A	C5-C6-N1	-5.26	115.07	117.70
53	B5	2720	G	C5-C6-O6	-5.26	125.44	128.60
53	B5	2779	A	C5-C6-N6	-5.26	119.49	123.70
53	B5	3245	A	C5-C6-N6	-5.26	119.49	123.70
53	B5	3342	A	O4'-C1'-N9	5.26	112.41	108.20
1	AA	53	G	O4'-C1'-N9	5.26	112.41	108.20
1	AA	812	A	O4'-C1'-N9	5.26	112.41	108.20
1	AA	1386	C	O4'-C1'-N1	5.26	112.41	108.20
53	B5	2414	G	O4'-C1'-N9	5.26	112.41	108.20
1	AA	1384	G	C5-C6-O6	-5.26	125.45	128.60
1	AA	1424	A	C5-C6-N6	-5.26	119.49	123.70
1	AA	1519	G	OP1-P-OP2	-5.26	111.72	119.60
9	AJ	90	TYR	CB-CG-CD2	-5.26	117.85	121.00
29	BE	142	PHE	CB-CG-CD2	-5.26	117.12	120.80
53	B5	313	A	O4'-C1'-N9	5.26	112.41	108.20
53	B5	1423	C	N3-C4-N4	5.26	121.68	118.00
53	B5	1858	A	C4-C5-C6	5.26	119.63	117.00
53	B5	1882	G	O4'-C1'-N9	5.26	112.41	108.20
53	B5	2119	A	O4'-C1'-N9	5.26	112.40	108.20
1	AA	93	A	O4'-C1'-N9	5.25	112.40	108.20
1	AA	571	G	C5-C6-O6	-5.25	125.45	128.60
53	B5	3227	A	O4'-C1'-N9	5.25	112.40	108.20
1	AA	91	G	O4'-C1'-N9	5.25	112.40	108.20
1	AA	245	U	O4'-C1'-N1	5.25	112.40	108.20
1	AA	550	A	O4'-C1'-N9	5.25	112.40	108.20
1	AA	1300	U	O4'-C1'-N1	5.25	112.40	108.20
1	AA	1456	G	O4'-C1'-N9	5.25	112.40	108.20
53	B5	1003	A	C5-C6-N6	-5.25	119.50	123.70
53	B5	3026	G	C5-C6-O6	-5.25	125.45	128.60
1	AA	480	G	O4'-C1'-N9	5.25	112.40	108.20
19	A7	35	A	N9-C1'-C2'	-5.25	106.22	112.00
53	B5	67	A	C5-C6-N6	-5.25	119.50	123.70
53	B5	934	G	O4'-C1'-N9	5.25	112.40	108.20
53	B5	2302	G	O4'-C1'-N9	5.25	112.40	108.20
1	AA	804	A	C5-C6-N6	-5.25	119.50	123.70
1	AA	1669	A	O4'-C1'-N9	5.25	112.40	108.20
1	AA	1778	G	O4'-C1'-N9	5.25	112.40	108.20
53	B5	815	G	O4'-C1'-N9	5.25	112.40	108.20
53	B5	1005	G	O4'-C1'-N9	5.25	112.40	108.20
53	B5	2817	A	C4-C5-C6	5.25	119.62	117.00
53	B5	2990	G	C5-C6-O6	-5.25	125.45	128.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	580	A	O4'-C1'-N9	5.25	112.40	108.20
53	B5	330	G	O4'-C1'-N9	5.25	112.40	108.20
53	B5	348	A	O4'-C1'-N9	5.25	112.40	108.20
53	B5	375	A	C4-C5-C6	5.25	119.62	117.00
53	B5	1159	A	O4'-C1'-N9	5.25	112.40	108.20
53	B5	1758	G	C5-C6-O6	-5.25	125.45	128.60
53	B5	1933	A	O4'-C1'-N9	5.25	112.40	108.20
53	B5	2415	C	N3-C4-N4	5.25	121.67	118.00
53	B5	2467	G	O4'-C1'-N9	5.25	112.40	108.20
53	B5	3336	A	C5-C6-N6	-5.25	119.50	123.70
1	AA	914	A	C4-C5-C6	5.25	119.62	117.00
1	AA	1239	G	O4'-C1'-N9	5.25	112.40	108.20
53	B5	577	C	O4'-C1'-N1	5.25	112.40	108.20
53	B5	2597	U	O4'-C1'-N1	5.25	112.40	108.20
53	B5	2642	A	C4-C5-C6	5.25	119.62	117.00
1	AA	927	U	O4'-C1'-N1	5.25	112.40	108.20
1	AA	929	A	O4'-C1'-N9	5.25	112.40	108.20
53	B5	1934	G	N1-C6-O6	5.25	123.05	119.90
1	AA	204	G	N1-C6-O6	5.24	123.05	119.90
1	AA	1598	A	C4-C5-C6	5.24	119.62	117.00
53	B5	320	G	C5-C6-O6	-5.24	125.45	128.60
53	B5	789	A	O4'-C1'-N9	5.24	112.40	108.20
53	B5	991	U	O4'-C1'-N1	5.24	112.39	108.20
53	B5	1286	A	O4'-C1'-N9	5.24	112.39	108.20
53	B5	2257	C	O4'-C1'-N1	5.24	112.39	108.20
53	B5	2474	G	C5-C6-O6	-5.24	125.45	128.60
53	B5	3372	A	C4-C5-C6	5.24	119.62	117.00
1	AA	219	A	C5-C6-N6	-5.24	119.51	123.70
8	AI	126	PRO	C-N-CA	5.24	134.81	121.70
53	B5	820	A	C4-C5-C6	5.24	119.62	117.00
53	B5	1909	A	O4'-C1'-N9	5.24	112.39	108.20
53	B5	3186	A	O4'-C1'-N9	5.24	112.39	108.20
1	AA	417	A	C5-C6-N6	-5.24	119.51	123.70
1	AA	737	A	O4'-C1'-N9	5.24	112.39	108.20
1	AA	1439	U	O4'-C1'-N1	5.24	112.39	108.20
1	AA	1695	G	N1-C6-O6	5.24	123.04	119.90
51	B3	49	G	C5-C6-O6	-5.24	125.46	128.60
53	B5	668	G	O4'-C1'-N9	5.24	112.39	108.20
53	B5	1110	U	O4'-C1'-N1	5.24	112.39	108.20
53	B5	2396	G	C5-C6-O6	-5.24	125.46	128.60
53	B5	3168	A	C5-C6-N6	-5.24	119.51	123.70
1	AA	202	A	C5-C6-N6	-5.24	119.51	123.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	858	G	C4-N9-C1'	5.24	133.31	126.50
29	BE	22	ARG	NE-CZ-NH2	-5.24	117.68	120.30
53	B5	288	C	N3-C4-N4	5.24	121.67	118.00
53	B5	338	A	C4-C5-C6	5.24	119.62	117.00
53	B5	390	G	O4'-C1'-N9	5.24	112.39	108.20
53	B5	968	G	O4'-C1'-N9	5.24	112.39	108.20
53	B5	1202	A	C4-C5-C6	5.24	119.62	117.00
53	B5	1580	A	C4-C5-C6	5.24	119.62	117.00
53	B5	1889	G	C5-C6-O6	-5.24	125.46	128.60
53	B5	2172	A	P-O3'-C3'	5.24	125.99	119.70
1	AA	1521	G	P-O5'-C5'	-5.24	112.52	120.90
1	AA	505	A	C5-C6-N6	-5.24	119.51	123.70
53	B5	665	A	O4'-C1'-N9	5.24	112.39	108.20
53	B5	1776	G	O4'-C1'-N9	5.24	112.39	108.20
53	B5	1900	A	C4-C5-C6	5.24	119.62	117.00
53	B5	1921	A	C5-C6-N1	-5.24	115.08	117.70
53	B5	2113	A	C5-C6-N6	-5.24	119.51	123.70
53	B5	2562	G	O4'-C1'-N9	5.24	112.39	108.20
1	AA	164	A	O4'-C1'-N9	5.23	112.39	108.20
1	AA	707	A	C5-C6-N6	-5.23	119.51	123.70
1	AA	1668	G	C5-C6-O6	-5.23	125.46	128.60
53	B5	1756	C	N3-C4-N4	5.23	121.66	118.00
53	B5	2888	U	O4'-C1'-N1	5.23	112.39	108.20
1	AA	1184	G	C5-C6-O6	-5.23	125.46	128.60
1	AA	1419	A	C4-C5-C6	5.23	119.62	117.00
53	B5	183	G	O4'-C1'-N9	5.23	112.39	108.20
53	B5	523	A	O4'-C1'-N9	5.23	112.39	108.20
53	B5	614	C	N3-C4-N4	5.23	121.66	118.00
53	B5	1002	A	O4'-C1'-N9	5.23	112.39	108.20
53	B5	2198	A	P-O3'-C3'	5.23	125.98	119.70
53	B5	2458	A	O4'-C1'-N9	5.23	112.39	108.20
53	B5	2711	C	N3-C4-N4	5.23	121.66	118.00
53	B5	3183	A	O4'-C1'-N9	5.23	112.38	108.20
1	AA	1030	U	O4'-C1'-N1	5.23	112.38	108.20
52	B4	75	G	C5-C6-O6	-5.23	125.46	128.60
53	B5	2155	G	C5-C6-O6	-5.23	125.46	128.60
53	B5	3238	G	C5-C6-O6	-5.23	125.46	128.60
1	AA	685	A	O4'-C1'-N9	5.23	112.38	108.20
1	AA	687	G	N1-C6-O6	5.23	123.04	119.90
1	AA	966	A	C5-C6-N6	-5.23	119.52	123.70
1	AA	1024	A	O4'-C1'-N9	5.23	112.38	108.20
1	AA	1223	A	O4'-C1'-N9	5.23	112.38	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	417	A	C5-C6-N6	-5.23	119.52	123.70
53	B5	1203	A	O4'-C1'-N9	5.23	112.38	108.20
53	B5	2418	G	C5-C6-O6	-5.23	125.46	128.60
53	B5	2811	A	C4-C5-C6	5.23	119.61	117.00
53	B5	3202	G	O4'-C1'-N9	5.23	112.38	108.20
1	AA	440	U	O4'-C1'-N1	5.23	112.38	108.20
53	B5	1697	A	C5-C6-N6	-5.23	119.52	123.70
53	B5	2837	A	C4-C5-C6	5.23	119.61	117.00
1	AA	80	A	O4'-C1'-N9	5.22	112.38	108.20
1	AA	1173	G	O4'-C1'-N9	5.22	112.38	108.20
41	BQ	14	MET	CG-SD-CE	-5.22	91.84	100.20
53	B5	402	A	C4-C5-C6	5.22	119.61	117.00
53	B5	1491	A	C4-C5-C6	5.22	119.61	117.00
53	B5	1603	A	C4-C5-C6	5.22	119.61	117.00
53	B5	2325	G	O4'-C1'-N9	5.22	112.38	108.20
53	B5	2452	G	O4'-C1'-N9	5.22	112.38	108.20
53	B5	3126	C	C5'-C4'-O4'	5.22	115.37	109.10
53	B5	3200	G	O4'-C1'-N9	5.22	112.38	108.20
53	B5	3322	A	C4-C5-C6	5.22	119.61	117.00
1	AA	707	A	O4'-C1'-N9	5.22	112.38	108.20
1	AA	1521	G	C4-N9-C1'	5.22	133.29	126.50
1	AA	1548	A	C1'-O4'-C4'	-5.22	105.72	109.90
52	B4	92	A	O4'-C1'-N9	5.22	112.38	108.20
53	B5	967	A	O4'-C1'-N9	5.22	112.38	108.20
53	B5	995	G	C5-C6-O6	-5.22	125.47	128.60
53	B5	1595	U	P-O3'-C3'	5.22	125.97	119.70
53	B5	2667	A	O4'-C1'-N9	5.22	112.38	108.20
1	AA	1303	C	N3-C4-N4	5.22	121.66	118.00
19	A7	70	C	N3-C4-N4	-5.22	114.34	118.00
53	B5	920	A	C4-C5-C6	5.22	119.61	117.00
53	B5	1245	A	C4-C5-C6	5.22	119.61	117.00
53	B5	2925	C	N3-C4-N4	5.22	121.66	118.00
53	B5	3099	C	C2-N1-C1'	5.22	124.54	118.80
1	AA	823	G	C5-C6-O6	-5.22	125.47	128.60
1	AA	937	G	O4'-C1'-N9	5.22	112.38	108.20
1	AA	1146	G	C5-C6-O6	-5.22	125.47	128.60
1	AA	1380	G	C5-C6-O6	-5.22	125.47	128.60
12	AM	85	PHE	CB-CG-CD2	-5.22	117.15	120.80
52	B4	140	G	C5-C6-O6	-5.22	125.47	128.60
53	B5	256	G	C5-C6-O6	-5.22	125.47	128.60
53	B5	917	A	C4-C5-C6	5.22	119.61	117.00
53	B5	2215	A	C5-C6-N6	-5.22	119.53	123.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	2527	G	C5-C6-O6	-5.22	125.47	128.60
53	B5	2926	A	C4-C5-C6	5.22	119.61	117.00
53	B5	3170	A	C4-C5-C6	5.22	119.61	117.00
1	AA	906	A	O4'-C1'-N9	5.22	112.37	108.20
19	A7	18	G	C5-C6-O6	5.22	131.73	128.60
19	A7	51	G	C4-C5-N7	-5.22	108.71	110.80
53	B5	632	G	C5-C6-O6	-5.22	125.47	128.60
53	B5	1085	A	C5-C6-N6	-5.22	119.53	123.70
53	B5	1647	A	C4-C5-C6	5.22	119.61	117.00
1	AA	344	A	O4'-C1'-N9	5.22	112.37	108.20
1	AA	780	A	O4'-C1'-N9	5.22	112.37	108.20
1	AA	1073	A	C4-C5-C6	5.22	119.61	117.00
1	AA	1160	A	C5-C6-N1	-5.22	115.09	117.70
19	A7	66	A	C3'-C2'-C1'	5.22	105.67	101.50
38	BN	60	PHE	CB-CG-CD2	-5.22	117.15	120.80
53	B5	259	C	N3-C4-C5	-5.22	119.81	121.90
53	B5	841	A	O4'-C1'-N9	5.22	112.37	108.20
53	B5	925	A	C4-C5-C6	5.22	119.61	117.00
53	B5	1106	G	C5-C6-O6	-5.22	125.47	128.60
53	B5	1176	C	O4'-C1'-N1	5.22	112.37	108.20
53	B5	2564	G	O4'-C1'-N9	5.22	112.37	108.20
53	B5	2769	A	C5-C6-N1	-5.22	115.09	117.70
53	B5	189	G	P-O3'-C3'	5.21	125.96	119.70
53	B5	965	A	C4-C5-C6	5.21	119.61	117.00
53	B5	1418	A	O4'-C1'-N9	5.21	112.37	108.20
53	B5	2733	A	O4'-C1'-N9	5.21	112.37	108.20
53	B5	2752	U	O4'-C1'-N1	5.21	112.37	108.20
53	B5	2993	G	O4'-C1'-N9	5.21	112.37	108.20
53	B5	3391	A	O4'-C1'-N9	5.21	112.37	108.20
1	AA	1084	A	O4'-C1'-N9	5.21	112.37	108.20
51	B3	112	G	C5-C6-O6	-5.21	125.47	128.60
1	AA	1155	C	N3-C4-N4	5.21	121.65	118.00
1	AA	1761	A	C4-C5-C6	5.21	119.61	117.00
51	B3	92	A	O4'-C1'-N9	5.21	112.37	108.20
52	B4	120	C	N3-C4-N4	5.21	121.65	118.00
53	B5	1532	C	N3-C4-N4	5.21	121.65	118.00
53	B5	1805	C	O4'-C1'-N1	5.21	112.37	108.20
53	B5	1893	A	C5-C6-N6	-5.21	119.53	123.70
53	B5	2256	A	C5-C6-N6	-5.21	119.53	123.70
1	AA	67	A	O4'-C1'-N9	5.21	112.37	108.20
53	B5	39	A	C5-C6-N1	-5.21	115.09	117.70
53	B5	2770	G	O4'-C1'-N9	5.21	112.37	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	189	C	N3-C4-N4	5.21	121.65	118.00
1	AA	814	A	O4'-C1'-N9	5.21	112.37	108.20
1	AA	1233	G	C5-C6-O6	-5.21	125.47	128.60
1	AA	1716	G	C5-C6-O6	-5.21	125.47	128.60
19	A7	67	A	C5-C6-N6	5.21	127.87	123.70
53	B5	222	A	C5-C6-N6	-5.21	119.53	123.70
53	B5	783	A	C4-C5-C6	5.21	119.61	117.00
53	B5	1521	G	O4'-C1'-N9	5.21	112.37	108.20
53	B5	1538	G	O4'-C1'-N9	5.21	112.37	108.20
1	AA	264	G	O4'-C1'-N9	5.21	112.36	108.20
1	AA	576	G	O4'-C1'-N9	5.21	112.37	108.20
1	AA	1520	U	C5'-C4'-C3'	-5.21	107.67	116.00
53	B5	18	G	C5-C6-O6	-5.21	125.48	128.60
53	B5	809	G	O4'-C1'-N9	5.21	112.36	108.20
53	B5	1126	G	C5-C6-O6	-5.21	125.48	128.60
53	B5	1426	C	N3-C4-C5	-5.21	119.82	121.90
53	B5	2175	U	O4'-C1'-N1	5.21	112.37	108.20
1	AA	68	A	C4-C5-C6	5.21	119.60	117.00
1	AA	1717	A	O4'-C1'-N9	5.20	112.36	108.20
53	B5	51	A	O4'-C1'-N9	5.20	112.36	108.20
53	B5	383	G	O4'-C1'-N9	5.20	112.36	108.20
53	B5	588	G	C5-C6-O6	-5.20	125.48	128.60
53	B5	1075	A	C5-C6-N6	-5.20	119.54	123.70
53	B5	1268	G	C5-C6-O6	-5.20	125.48	128.60
53	B5	1618	G	O4'-C1'-N9	5.20	112.36	108.20
53	B5	2442	G	C5-C6-O6	-5.20	125.48	128.60
53	B5	2539	A	C5-C6-N6	-5.20	119.54	123.70
53	B5	3193	C	N3-C4-C5	-5.20	119.82	121.90
1	AA	1028	U	O4'-C1'-N1	5.20	112.36	108.20
1	AA	1503	A	O4'-C1'-N9	5.20	112.36	108.20
8	AI	138	PHE	CB-CG-CD2	5.20	124.44	120.80
19	A7	73	A	N9-C1'-C2'	-5.20	106.28	112.00
53	B5	1841	A	O4'-C1'-N9	5.20	112.36	108.20
19	A7	25	C	C3'-C2'-C1'	-5.20	97.34	101.50
19	A7	57	G	N3-C2-N2	-5.20	116.26	119.90
53	B5	172	G	N1-C6-O6	5.20	123.02	119.90
53	B5	791	A	C4-C5-C6	5.20	119.60	117.00
53	B5	830	A	C5-C6-N6	-5.20	119.54	123.70
53	B5	1254	C	N3-C4-N4	5.20	121.64	118.00
53	B5	1754	G	C5-C6-O6	-5.20	125.48	128.60
53	B5	2676	A	C4-C5-C6	5.20	119.60	117.00
53	B5	2697	A	O4'-C1'-N9	5.20	112.36	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	3073	A	C5-C6-N1	-5.20	115.10	117.70
1	AA	1588	G	O4'-C1'-N9	5.20	112.36	108.20
53	B5	80	G	C5-C6-O6	-5.20	125.48	128.60
53	B5	1370	G	C5-C6-O6	-5.20	125.48	128.60
53	B5	2903	A	C5-C6-N6	-5.20	119.54	123.70
53	B5	3029	A	O4'-C1'-N9	5.20	112.36	108.20
53	B5	2290	C	N3-C4-C5	-5.20	119.82	121.90
1	AA	240	U	C6-N1-C1'	-5.20	113.92	121.20
1	AA	336	G	C5-C6-O6	-5.20	125.48	128.60
1	AA	446	A	O4'-C1'-N9	5.20	112.36	108.20
53	B5	1149	G	P-O3'-C3'	5.20	125.93	119.70
53	B5	1787	A	C4-C5-C6	5.20	119.60	117.00
53	B5	2929	C	N3-C4-N4	5.20	121.64	118.00
1	AA	401	A	C4-C5-C6	5.19	119.60	117.00
53	B5	39	A	O4'-C1'-N9	5.19	112.36	108.20
53	B5	651	G	P-O3'-C3'	-5.19	113.47	119.70
53	B5	2489	C	N3-C4-C5	-5.19	119.82	121.90
53	B5	2745	G	C5-C6-O6	-5.19	125.48	128.60
53	B5	2801	A	C4-C5-C6	5.19	119.60	117.00
1	AA	1273	G	P-O3'-C3'	5.19	125.93	119.70
1	AA	1666	G	C5-C6-O6	-5.19	125.48	128.60
53	B5	61	A	P-O3'-C3'	5.19	125.93	119.70
53	B5	1132	C	N3-C4-N4	5.19	121.63	118.00
53	B5	1417	G	C5-C6-O6	-5.19	125.48	128.60
53	B5	1759	C	N3-C4-N4	5.19	121.64	118.00
53	B5	2367	A	C5-C6-N6	-5.19	119.55	123.70
53	B5	2674	A	O4'-C1'-N9	5.19	112.35	108.20
53	B5	3362	A	O4'-C1'-N9	5.19	112.35	108.20
1	AA	82	U	O4'-C1'-N1	5.19	112.35	108.20
1	AA	386	G	O4'-C1'-N9	5.19	112.35	108.20
1	AA	407	A	C5-C6-N1	-5.19	115.11	117.70
1	AA	1217	A	O4'-C1'-N9	5.19	112.35	108.20
1	AA	1592	G	C5-C6-O6	-5.19	125.49	128.60
19	A7	56	C	N3-C4-C5	-5.19	119.82	121.90
51	B3	34	C	N3-C4-N4	5.19	121.63	118.00
53	B5	247	C	N3-C4-C5	-5.19	119.82	121.90
53	B5	1145	G	N1-C6-O6	5.19	123.01	119.90
53	B5	1155	C	N3-C4-C5	-5.19	119.82	121.90
53	B5	1504	A	C4-C5-C6	5.19	119.59	117.00
53	B5	2116	G	O4'-C1'-N9	5.19	112.35	108.20
53	B5	2152	A	O4'-C1'-N9	5.19	112.35	108.20
53	B5	1793	C	N3-C4-N4	5.19	121.63	118.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	2559	A	C4-C5-C6	5.19	119.59	117.00
1	AA	105	A	P-O3'-C3'	5.19	125.92	119.70
1	AA	187	G	O4'-C1'-N9	5.19	112.35	108.20
53	B5	1647	A	O4'-C1'-N9	5.19	112.35	108.20
53	B5	2813	A	O4'-C1'-N9	5.19	112.35	108.20
1	AA	1285	G	C5-C6-O6	-5.19	125.49	128.60
19	A7	74	C	C5-C6-N1	-5.19	118.41	121.00
1	AA	417	A	C4-C5-C6	5.18	119.59	117.00
1	AA	938	A	C4-C5-C6	5.18	119.59	117.00
1	AA	1398	A	C4-C5-C6	5.18	119.59	117.00
19	A7	3	G	C6-C5-N7	-5.18	127.29	130.40
19	A7	69	U	C4-C5-C6	5.18	122.81	119.70
24	B9	36	ARG	NE-CZ-NH1	5.18	122.89	120.30
52	B4	89	A	C5-C6-N6	-5.18	119.55	123.70
53	B5	77	A	O4'-C1'-N9	5.18	112.35	108.20
53	B5	100	A	C5-C6-N6	-5.18	119.55	123.70
53	B5	621	A	O4'-C1'-N9	5.18	112.35	108.20
53	B5	2253	G	C5-C6-O6	-5.18	125.49	128.60
53	B5	2864	A	C5'-C4'-O4'	5.18	115.32	109.10
53	B5	3021	A	C5-C6-N1	-5.18	115.11	117.70
1	AA	301	A	C4-C5-C6	5.18	119.59	117.00
1	AA	901	G	O4'-C1'-N9	5.18	112.34	108.20
53	B5	409	A	C5-C6-N6	-5.18	119.55	123.70
53	B5	432	G	O4'-C1'-N9	5.18	112.35	108.20
53	B5	1842	A	C4-C5-C6	5.18	119.59	117.00
53	B5	2839	G	O4'-C1'-N9	5.18	112.35	108.20
1	AA	468	A	C4-C5-C6	5.18	119.59	117.00
1	AA	1130	A	C4-C5-C6	5.18	119.59	117.00
53	B5	193	C	N3-C4-C5	-5.18	119.83	121.90
53	B5	940	G	O4'-C1'-N9	5.18	112.34	108.20
53	B5	1915	A	C5-C6-N1	-5.18	115.11	117.70
53	B5	2527	G	O4'-C1'-N9	5.18	112.34	108.20
1	AA	68	A	C5-C6-N6	-5.18	119.56	123.70
1	AA	1415	G	O4'-C1'-N9	5.18	112.34	108.20
1	AA	1416	G	O4'-C1'-N9	5.18	112.34	108.20
1	AA	1422	A	C5-C6-N6	-5.18	119.56	123.70
52	B4	52	A	C4-C5-C6	5.18	119.59	117.00
53	B5	102	C	N3-C4-N4	5.18	121.62	118.00
53	B5	787	G	O4'-C1'-N9	5.18	112.34	108.20
53	B5	1624	G	O4'-C1'-N9	5.18	112.34	108.20
52	B4	48	A	O4'-C1'-N9	5.18	112.34	108.20
1	AA	151	G	C5-C6-O6	-5.18	125.49	128.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	928	A	C5-C6-N6	-5.18	119.56	123.70
52	B4	37	A	O4'-C1'-N9	5.18	112.34	108.20
52	B4	110	C	N3-C4-C5	-5.18	119.83	121.90
53	B5	361	A	O4'-C1'-N9	5.18	112.34	108.20
53	B5	2540	A	C5-C6-N6	-5.18	119.56	123.70
53	B5	2691	A	O4'-C1'-N9	5.18	112.34	108.20
1	AA	1581	A	C4-C5-C6	5.17	119.59	117.00
53	B5	217	U	P-O3'-C3'	5.17	125.91	119.70
53	B5	573	C	N3-C4-C5	-5.17	119.83	121.90
53	B5	639	G	O4'-C1'-N9	5.17	112.34	108.20
53	B5	2961	G	O4'-C1'-N9	5.17	112.34	108.20
1	AA	671	G	C5-C6-O6	-5.17	125.50	128.60
1	AA	1226	A	O4'-C1'-N9	5.17	112.34	108.20
53	B5	194	U	O4'-C1'-N1	5.17	112.34	108.20
53	B5	1319	G	O4'-C1'-N9	5.17	112.34	108.20
1	AA	551	G	C5-C6-O6	-5.17	125.50	128.60
1	AA	820	U	O4'-C1'-N1	5.17	112.34	108.20
1	AA	1136	A	C4-C5-C6	5.17	119.59	117.00
1	AA	1730	A	O4'-C1'-N9	5.17	112.34	108.20
52	B4	43	A	C5-C6-N6	-5.17	119.56	123.70
53	B5	1469	C	C6-N1-C1'	-5.17	114.59	120.80
53	B5	2122	G	N1-C6-O6	5.17	123.00	119.90
53	B5	2330	C	N3-C4-C5	-5.17	119.83	121.90
1	AA	660	G	O4'-C1'-N9	5.17	112.34	108.20
19	A7	14	A	C6-N1-C2	-5.17	115.50	118.60
53	B5	806	A	C4-C5-C6	5.17	119.58	117.00
53	B5	2120	A	C4-C5-C6	5.17	119.58	117.00
1	AA	496	G	C5-C6-O6	-5.17	125.50	128.60
19	A7	73	A	C4-C5-C6	-5.17	114.42	117.00
53	B5	202	G	N1-C6-O6	5.17	123.00	119.90
53	B5	750	G	O4'-C1'-N9	5.17	112.33	108.20
53	B5	2419	A	C5-C6-N6	-5.17	119.56	123.70
53	B5	2682	C	N3-C4-N4	5.17	121.62	118.00
53	B5	2750	U	O4'-C1'-N1	5.17	112.33	108.20
1	AA	19	A	O4'-C1'-N9	5.17	112.33	108.20
1	AA	571	G	O4'-C1'-N9	5.17	112.33	108.20
1	AA	895	G	C5-C6-O6	-5.17	125.50	128.60
1	AA	949	C	N3-C4-N4	5.17	121.62	118.00
1	AA	1653	A	C4-C5-C6	5.17	119.58	117.00
51	B3	86	G	C5-C6-O6	-5.17	125.50	128.60
52	B4	40	A	C5-C6-N6	-5.17	119.57	123.70
52	B4	69	U	P-O5'-C5'	-5.17	112.63	120.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	12	A	C5-C6-N6	-5.17	119.57	123.70
53	B5	1787	A	O4'-C1'-N9	5.17	112.33	108.20
53	B5	1915	A	O4'-C1'-N9	5.17	112.33	108.20
53	B5	3004	C	N3-C4-C5	-5.17	119.83	121.90
53	B5	3366	G	O4'-C1'-N9	5.17	112.33	108.20
1	AA	1630	C	O5'-P-OP1	5.16	116.90	110.70
53	B5	100	A	O4'-C1'-N9	5.16	112.33	108.20
53	B5	910	G	O4'-C1'-N9	5.16	112.33	108.20
53	B5	1383	G	O4'-C1'-N9	5.16	112.33	108.20
53	B5	1390	A	O4'-C1'-N9	5.16	112.33	108.20
53	B5	2256	A	O4'-C1'-N9	5.16	112.33	108.20
53	B5	2649	A	O4'-C1'-N9	5.16	112.33	108.20
53	B5	2958	A	C5-C6-N6	-5.16	119.57	123.70
53	B5	1757	A	C4-C5-C6	5.16	119.58	117.00
1	AA	95	G	C5-C6-O6	-5.16	125.50	128.60
1	AA	1071	G	O4'-C1'-N9	5.16	112.33	108.20
52	B4	24	G	O4'-C1'-N9	5.16	112.33	108.20
52	B4	123	G	C5-C6-O6	-5.16	125.50	128.60
53	B5	763	G	C5-C6-O6	-5.16	125.50	128.60
53	B5	1203	A	C4-C5-C6	5.16	119.58	117.00
53	B5	1496	C	C2-N1-C1'	5.16	124.48	118.80
53	B5	2178	A	C4-C5-C6	5.16	119.58	117.00
53	B5	2219	A	C4-C5-C6	5.16	119.58	117.00
53	B5	2357	A	C5-C6-N6	-5.16	119.57	123.70
53	B5	2490	C	N3-C4-N4	5.16	121.61	118.00
1	AA	109	G	O4'-C1'-N9	5.16	112.33	108.20
1	AA	677	G	C5-C6-O6	-5.16	125.50	128.60
53	B5	285	A	C4-C5-C6	5.16	119.58	117.00
53	B5	418	A	C5-C6-N6	-5.16	119.57	123.70
53	B5	1776	G	P-O3'-C3'	5.16	125.89	119.70
52	B4	14	C	N3-C4-C5	-5.16	119.84	121.90
53	B5	2637	A	O4'-C1'-N9	5.16	112.33	108.20
1	AA	23	G	C5-C6-O6	-5.16	125.51	128.60
1	AA	273	G	O4'-C1'-N9	5.16	112.32	108.20
1	AA	962	A	O4'-C1'-N9	5.16	112.32	108.20
1	AA	1451	G	O4'-C1'-N9	5.16	112.33	108.20
53	B5	1013	G	O4'-C1'-N9	5.16	112.32	108.20
53	B5	1514	G	C5-C6-O6	-5.16	125.51	128.60
53	B5	1901	A	C4-C5-C6	5.16	119.58	117.00
53	B5	2353	G	O4'-C1'-N9	5.16	112.33	108.20
53	B5	2424	A	O4'-C1'-N9	5.16	112.32	108.20
53	B5	2640	A	C4-C5-C6	5.16	119.58	117.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	3275	A	O4'-C1'-N9	5.16	112.32	108.20
1	AA	1761	A	C5-C6-N6	-5.15	119.58	123.70
53	B5	70	A	C5-C6-N6	-5.15	119.58	123.70
53	B5	157	A	C4-C5-C6	5.15	119.58	117.00
53	B5	593	A	O4'-C1'-N9	5.15	112.32	108.20
53	B5	695	C	N3-C4-C5	-5.15	119.84	121.90
53	B5	1704	A	C5-C6-N6	-5.15	119.58	123.70
53	B5	1787	A	C5-C6-N6	-5.15	119.58	123.70
1	AA	730	G	O4'-C1'-N9	5.15	112.32	108.20
1	AA	1217	A	C5-C6-N6	-5.15	119.58	123.70
53	B5	131	C	N3-C4-N4	5.15	121.61	118.00
53	B5	503	C	N3-C4-N4	5.15	121.61	118.00
53	B5	618	C	N3-C4-C5	-5.15	119.84	121.90
53	B5	621	A	C4-C5-C6	5.15	119.58	117.00
53	B5	697	A	O4'-C1'-N9	5.15	112.32	108.20
53	B5	1676	A	C5-C6-N6	-5.15	119.58	123.70
53	B5	3264	G	O4'-C1'-N9	5.15	112.32	108.20
1	AA	85	A	O4'-C1'-N9	5.15	112.32	108.20
1	AA	1006	C	N3-C4-N4	5.15	121.61	118.00
1	AA	1260	G	C5-C6-O6	-5.15	125.51	128.60
51	B3	46	A	C4-C5-C6	5.15	119.58	117.00
53	B5	1285	G	O4'-C1'-N9	5.15	112.32	108.20
53	B5	2947	G	C5-C6-O6	-5.15	125.51	128.60
1	AA	977	A	O4'-C1'-N9	5.15	112.32	108.20
1	AA	1484	G	C5-C6-O6	-5.15	125.51	128.60
53	B5	1454	A	C4-C5-C6	5.15	119.57	117.00
53	B5	3253	G	C5-C6-O6	-5.15	125.51	128.60
53	B5	810	A	O4'-C1'-N9	5.15	112.32	108.20
53	B5	1450	G	O4'-C1'-N9	5.15	112.32	108.20
53	B5	1872	C	N3-C4-C5	-5.15	119.84	121.90
53	B5	2398	A	C4-C5-C6	5.15	119.57	117.00
53	B5	2740	A	C4-C5-C6	5.15	119.57	117.00
1	AA	864	U	O3'-P-O5'	5.14	113.78	104.00
1	AA	903	G	C5-C6-O6	-5.14	125.51	128.60
1	AA	1475	G	O4'-C1'-N9	5.14	112.32	108.20
19	A7	69	U	C5-C6-N1	-5.14	120.13	122.70
53	B5	1825	G	O4'-C1'-N9	5.14	112.31	108.20
53	B5	1847	A	C4-C5-C6	5.14	119.57	117.00
53	B5	1918	C	N3-C4-C5	-5.14	119.84	121.90
53	B5	2689	A	C4-C5-C6	5.14	119.57	117.00
53	B5	3148	U	P-O3'-C3'	5.14	125.87	119.70
1	AA	529	A	O4'-C1'-N9	5.14	112.31	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	565	C	N3-C4-N4	5.14	121.60	118.00
53	B5	166	C	N3-C4-N4	5.14	121.60	118.00
53	B5	1120	A	O4'-C1'-N9	5.14	112.31	108.20
53	B5	1517	G	C5-C6-O6	-5.14	125.51	128.60
53	B5	2437	G	C5-C6-O6	-5.14	125.51	128.60
53	B5	2456	A	O4'-C1'-N9	5.14	112.31	108.20
53	B5	2736	A	C4-C5-C6	5.14	119.57	117.00
1	AA	412	A	O4'-C1'-N9	5.14	112.31	108.20
1	AA	1787	G	O4'-C1'-N9	5.14	112.31	108.20
52	B4	86	U	O4'-C1'-N1	5.14	112.31	108.20
53	B5	675	C	N3-C4-C5	-5.14	119.84	121.90
53	B5	1089	G	C5-C6-O6	-5.14	125.52	128.60
1	AA	1663	U	O4'-C1'-N1	5.14	112.31	108.20
1	AA	1792	A	C4-C5-C6	5.14	119.57	117.00
52	B4	67	U	O4'-C1'-N1	5.14	112.31	108.20
53	B5	961	C	O4'-C1'-N1	5.14	112.31	108.20
53	B5	2438	A	O4'-C1'-N9	5.14	112.31	108.20
53	B5	3076	C	N3-C4-N4	5.14	121.60	118.00
53	B5	3086	A	C5-C6-N1	-5.14	115.13	117.70
1	AA	1144	A	O4'-C1'-N9	5.14	112.31	108.20
52	B4	114	G	C5-C6-O6	-5.14	125.52	128.60
53	B5	718	G	C4-N9-C1'	5.14	133.18	126.50
53	B5	1013	G	C5-C6-O6	-5.14	125.52	128.60
53	B5	2448	G	N1-C6-O6	5.14	122.98	119.90
53	B5	2568	C	N3-C4-C5	-5.14	119.84	121.90
53	B5	2737	C	N3-C4-N4	5.14	121.60	118.00
53	B5	3140	G	O4'-C1'-N9	5.14	112.31	108.20
1	AA	933	C	O4'-C1'-N1	5.14	112.31	108.20
1	AA	940	A	O4'-C1'-N9	5.14	112.31	108.20
53	B5	1850	A	O4'-C1'-N9	5.14	112.31	108.20
53	B5	2371	G	O4'-C1'-N9	5.14	112.31	108.20
1	AA	1362	C	N3-C4-N4	5.13	121.59	118.00
53	B5	2958	A	C4-C5-C6	5.13	119.57	117.00
1	AA	484	C	N3-C4-N4	5.13	121.59	118.00
1	AA	1593	U	O4'-C1'-N1	5.13	112.31	108.20
1	AA	322	G	O4'-C1'-N9	5.13	112.31	108.20
1	AA	441	A	O4'-C1'-N9	5.13	112.31	108.20
1	AA	1089	A	C5-C6-N6	-5.13	119.59	123.70
1	AA	1462	G	C5-C6-O6	-5.13	125.52	128.60
1	AA	1676	A	O4'-C1'-N9	5.13	112.31	108.20
19	A7	1	G	C5'-C4'-C3'	5.13	124.21	116.00
53	B5	1843	C	N3-C4-C5	-5.13	119.85	121.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	2120	A	O4'-C1'-N9	5.13	112.31	108.20
53	B5	3271	A	O4'-C1'-N9	5.13	112.31	108.20
52	B4	9	A	C4-C5-C6	5.13	119.56	117.00
53	B5	3235	C	N3-C4-N4	5.13	121.59	118.00
1	AA	46	A	O4'-C1'-N9	5.13	112.30	108.20
1	AA	939	A	C4-C5-C6	5.13	119.56	117.00
1	AA	1354	A	C5-C6-N6	-5.13	119.60	123.70
1	AA	1730	A	C4-C5-C6	5.13	119.56	117.00
53	B5	365	A	C5-C6-N6	-5.13	119.60	123.70
53	B5	710	A	C4-C5-C6	5.13	119.56	117.00
53	B5	784	A	C5-C6-N1	-5.13	115.14	117.70
53	B5	951	A	O4'-C1'-N9	5.13	112.30	108.20
53	B5	2867	C	N3-C4-C5	-5.13	119.85	121.90
53	B5	2914	G	C5-C6-O6	-5.13	125.52	128.60
1	AA	86	A	O4'-C1'-N9	5.13	112.30	108.20
1	AA	1661	G	O4'-C1'-N9	5.13	112.30	108.20
53	B5	845	G	O4'-C1'-N9	5.13	112.30	108.20
53	B5	2580	A	O4'-C1'-N9	5.13	112.30	108.20
53	B5	2676	A	C5-C6-N6	-5.13	119.60	123.70
53	B5	2963	C	N3-C4-C5	-5.13	119.85	121.90
1	AA	370	A	C4-C5-C6	5.12	119.56	117.00
51	B3	85	G	C5-C6-O6	-5.12	125.53	128.60
1	AA	279	G	O4'-C1'-N9	5.12	112.30	108.20
1	AA	638	U	O4'-C1'-N1	5.12	112.30	108.20
8	AI	142	TYR	CA-CB-CG	-5.12	103.66	113.40
53	B5	832	G	O4'-C1'-N9	5.12	112.30	108.20
53	B5	1656	A	O4'-C1'-N9	5.12	112.30	108.20
53	B5	3126	C	N3-C4-C5	-5.12	119.85	121.90
1	AA	285	G	C5-C6-O6	-5.12	125.53	128.60
1	AA	1313	G	O4'-C1'-N9	5.12	112.30	108.20
1	AA	1464	G	O4'-C1'-N9	5.12	112.30	108.20
53	B5	1504	A	C5-C6-N6	-5.12	119.60	123.70
53	B5	1670	C	N3-C4-N4	5.12	121.58	118.00
53	B5	2369	G	C5-C6-O6	-5.12	125.53	128.60
53	B5	2479	C	N3-C4-N4	5.12	121.58	118.00
1	AA	646	C	N3-C4-N4	5.12	121.58	118.00
53	B5	826	G	O4'-C1'-N9	5.12	112.30	108.20
53	B5	1085	A	O4'-C1'-N9	5.12	112.30	108.20
53	B5	2555	G	C5-C6-O6	-5.12	125.53	128.60
53	B5	2956	A	C4-C5-C6	5.12	119.56	117.00
1	AA	178	U	O4'-C1'-N1	5.12	112.30	108.20
1	AA	534	A	O4'-C1'-N9	5.12	112.30	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1368	A	O4'-C1'-N9	5.12	112.30	108.20
1	AA	1473	A	O4'-C1'-N9	5.12	112.30	108.20
52	B4	144	G	P-O3'-C3'	5.12	125.84	119.70
53	B5	251	G	O4'-C1'-N9	5.12	112.29	108.20
53	B5	603	A	C4-C5-C6	5.12	119.56	117.00
53	B5	1528	G	O4'-C1'-N9	5.12	112.30	108.20
53	B5	2961	G	P-O3'-C3'	5.12	125.84	119.70
1	AA	1078	A	C4-C5-C6	5.12	119.56	117.00
53	B5	435	C	P-O3'-C3'	5.12	125.84	119.70
53	B5	975	G	C5-C6-O6	-5.12	125.53	128.60
53	B5	1741	A	C4-C5-C6	5.12	119.56	117.00
53	B5	2445	A	C5-C6-N6	-5.12	119.61	123.70
1	AA	1524	A	C5-C6-N6	-5.12	119.61	123.70
53	B5	798	G	C5-C6-O6	-5.12	125.53	128.60
53	B5	838	G	C5-C6-O6	-5.12	125.53	128.60
53	B5	1743	G	C5-C6-O6	-5.12	125.53	128.60
53	B5	2421	U	P-O3'-C3'	5.12	125.84	119.70
53	B5	2706	G	C5-C6-O6	-5.12	125.53	128.60
53	B5	2741	C	N3-C4-N4	5.12	121.58	118.00
53	B5	3021	A	C4-C5-C6	5.12	119.56	117.00
53	B5	3081	C	N3-C4-N4	5.12	121.58	118.00
53	B5	3086	A	C4-C5-C6	5.12	119.56	117.00
1	AA	42	G	C5-C6-O6	-5.11	125.53	128.60
1	AA	388	G	O4'-C1'-N9	5.11	112.29	108.20
1	AA	467	G	O4'-C1'-N9	5.11	112.29	108.20
1	AA	1385	A	C5-C6-N6	-5.11	119.61	123.70
1	AA	1568	A	O4'-C1'-N9	5.11	112.29	108.20
1	AA	1763	A	C4-C5-C6	5.11	119.56	117.00
53	B5	2628	A	C5-C6-N6	-5.11	119.61	123.70
1	AA	1683	G	C5-C6-O6	-5.11	125.53	128.60
53	B5	1529	A	C5-C6-N6	-5.11	119.61	123.70
53	B5	1868	G	O4'-C1'-N9	5.11	112.29	108.20
53	B5	3299	A	C5-C6-N6	-5.11	119.61	123.70
1	AA	425	A	C4-C5-C6	5.11	119.56	117.00
1	AA	778	G	C5-C6-O6	-5.11	125.53	128.60
53	B5	318	A	O4'-C1'-N9	5.11	112.29	108.20
53	B5	417	A	O4'-C1'-N9	5.11	112.29	108.20
53	B5	1830	G	O4'-C1'-N9	5.11	112.29	108.20
53	B5	1946	A	C4-C5-C6	5.11	119.56	117.00
1	AA	160	C	N3-C4-N4	5.11	121.58	118.00
1	AA	649	U	O4'-C1'-N1	5.11	112.29	108.20
53	B5	16	A	C4-C5-C6	5.11	119.55	117.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	1227	C	N3-C4-C5	-5.11	119.86	121.90
53	B5	1575	A	C5-C6-N6	-5.11	119.61	123.70
53	B5	1750	A	C4-C5-C6	5.11	119.55	117.00
53	B5	1883	A	C5-C6-N1	-5.11	115.14	117.70
53	B5	2360	C	O4'-C1'-N1	5.11	112.29	108.20
53	B5	3313	U	C5'-C4'-O4'	5.11	115.23	109.10
1	AA	490	C	N3-C4-N4	5.11	121.58	118.00
1	AA	500	C	N3-C4-C5	-5.11	119.86	121.90
1	AA	878	G	O4'-C1'-N9	5.11	112.29	108.20
1	AA	1001	G	O4'-C1'-N9	5.11	112.29	108.20
1	AA	1529	G	O4'-C1'-N9	5.11	112.29	108.20
53	B5	76	G	O4'-C1'-N9	5.11	112.29	108.20
53	B5	199	A	O4'-C1'-N9	5.11	112.29	108.20
53	B5	1332	A	O4'-C1'-N9	5.11	112.29	108.20
53	B5	1907	C	N3-C4-N4	5.11	121.58	118.00
53	B5	2733	A	C4-C5-C6	5.11	119.55	117.00
53	B5	3234	A	C4-C5-C6	5.11	119.55	117.00
51	B3	37	G	O4'-C1'-N9	5.11	112.28	108.20
53	B5	742	G	C5-C6-O6	-5.11	125.54	128.60
53	B5	796	U	O4'-C1'-N1	5.11	112.28	108.20
53	B5	1080	A	C5-C6-N6	-5.11	119.61	123.70
53	B5	1231	A	C5-C6-N6	-5.11	119.61	123.70
53	B5	2120	A	C5-C6-N6	-5.11	119.61	123.70
53	B5	2180	G	O4'-C1'-N9	5.11	112.28	108.20
53	B5	3218	A	O4'-C1'-N9	5.11	112.28	108.20
53	B5	3237	U	O4'-C1'-N1	5.11	112.28	108.20
53	B5	400	G	C5-C6-O6	-5.10	125.54	128.60
53	B5	2188	A	C5-C6-N6	-5.10	119.62	123.70
53	B5	2872	A	C4-C5-C6	5.10	119.55	117.00
1	AA	1133	U	C2'-C3'-O3'	5.10	121.86	113.70
52	B4	129	C	N3-C4-N4	5.10	121.57	118.00
53	B5	192	C	N3-C4-C5	-5.10	119.86	121.90
53	B5	1358	C	N3-C4-C5	-5.10	119.86	121.90
53	B5	2229	A	C5-C6-N6	-5.10	119.62	123.70
53	B5	2867	C	N3-C4-N4	5.10	121.57	118.00
1	AA	410	A	O4'-C1'-N9	5.10	112.28	108.20
1	AA	858	G	C4-C5-N7	5.10	112.84	110.80
8	AI	96	TYR	CB-CG-CD1	-5.10	117.94	121.00
19	A7	25	C	N3-C2-O2	-5.10	118.33	121.90
53	B5	45	A	O4'-C1'-N9	5.10	112.28	108.20
53	B5	547	G	O4'-C1'-N9	5.10	112.28	108.20
1	AA	329	G	C5-C6-O6	-5.10	125.54	128.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1467	A	O4'-C1'-N9	5.10	112.28	108.20
53	B5	151	A	C1'-O4'-C4'	-5.10	105.82	109.90
53	B5	355	A	C4-C5-C6	5.10	119.55	117.00
53	B5	368	G	C5-C6-O6	-5.10	125.54	128.60
53	B5	1066	G	C5-C6-O6	-5.10	125.54	128.60
53	B5	1574	C	N3-C4-N4	5.10	121.57	118.00
53	B5	1683	A	C4-C5-C6	5.10	119.55	117.00
53	B5	2825	C	N3-C4-N4	5.10	121.57	118.00
1	AA	438	A	C4-C5-C6	5.10	119.55	117.00
1	AA	1490	A	O4'-C1'-N9	5.10	112.28	108.20
1	AA	1537	G	N1-C6-O6	5.10	122.96	119.90
51	B3	101	A	C5-C6-N6	-5.10	119.62	123.70
53	B5	1726	C	N3-C4-N4	5.10	121.57	118.00
53	B5	1911	A	O4'-C1'-N9	5.10	112.28	108.20
53	B5	2494	A	O4'-C1'-N9	5.10	112.28	108.20
53	B5	535	G	O4'-C1'-N9	5.09	112.28	108.20
53	B5	1655	G	O4'-C1'-N9	5.09	112.28	108.20
53	B5	2710	C	N3-C4-N4	5.09	121.57	118.00
53	B5	3321	C	N3-C4-N4	5.09	121.57	118.00
1	AA	83	G	C5-C6-O6	-5.09	125.54	128.60
1	AA	1137	G	C5-C6-O6	-5.09	125.54	128.60
53	B5	2892	A	C5-C6-N6	-5.09	119.63	123.70
1	AA	67	A	C5-C6-N6	-5.09	119.63	123.70
1	AA	421	A	O4'-C1'-N9	5.09	112.27	108.20
1	AA	665	U	O4'-C1'-N1	5.09	112.27	108.20
1	AA	1333	A	O4'-C1'-N9	5.09	112.27	108.20
1	AA	1452	G	C5-C6-O6	-5.09	125.55	128.60
1	AA	1473	A	C4-C5-C6	5.09	119.55	117.00
53	B5	288	C	N3-C4-C5	-5.09	119.86	121.90
53	B5	402	A	C5-C6-N6	-5.09	119.63	123.70
53	B5	2146	C	N3-C4-N4	5.09	121.56	118.00
53	B5	2586	G	C5-C6-O6	-5.09	125.55	128.60
53	B5	2628	A	C4-C5-C6	5.09	119.55	117.00
53	B5	2869	U	O4'-C1'-N1	5.09	112.27	108.20
1	AA	96	G	C5-C6-O6	-5.09	125.55	128.60
1	AA	263	C	N3-C4-N4	5.09	121.56	118.00
1	AA	1026	A	C5-C6-N6	-5.09	119.63	123.70
1	AA	1692	A	C5-C6-N6	-5.09	119.63	123.70
19	A7	28	C	C1'-O4'-C4'	-5.09	105.83	109.90
53	B5	211	A	C4-C5-C6	5.09	119.55	117.00
53	B5	392	G	O4'-C1'-N9	5.09	112.27	108.20
53	B5	696	C	N3-C4-C5	-5.09	119.86	121.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	1335	C	N3-C4-N4	5.09	121.56	118.00
53	B5	2143	A	O4'-C1'-N9	5.09	112.27	108.20
53	B5	2802	A	C4-C5-C6	5.09	119.55	117.00
53	B5	2958	A	O4'-C1'-N9	5.09	112.27	108.20
1	AA	860	U	C1'-C2'-O2'	5.09	125.86	110.60
53	B5	805	G	O4'-C1'-N9	5.09	112.27	108.20
53	B5	1674	G	O4'-C1'-N9	5.09	112.27	108.20
53	B5	2143	A	C5-C6-N6	-5.09	119.63	123.70
1	AA	582	U	O4'-C1'-N1	5.09	112.27	108.20
1	AA	866	G	C1'-C2'-O2'	5.09	125.86	110.60
53	B5	154	U	O4'-C1'-N1	5.09	112.27	108.20
53	B5	552	G	P-O3'-C3'	5.09	125.80	119.70
53	B5	1566	A	O4'-C1'-N9	5.09	112.27	108.20
53	B5	1615	C	N3-C4-N4	5.09	121.56	118.00
53	B5	2138	A	O4'-C1'-N9	5.09	112.27	108.20
53	B5	2348	A	O4'-C1'-N9	5.09	112.27	108.20
53	B5	2413	A	O4'-C1'-N9	5.09	112.27	108.20
53	B5	2666	C	N3-C4-N4	5.09	121.56	118.00
53	B5	2887	A	C4-C5-C6	5.09	119.54	117.00
53	B5	2946	A	O4'-C1'-N9	5.09	112.27	108.20
53	B5	3062	G	O4'-C1'-N9	5.09	112.27	108.20
1	AA	521	A	O4'-C1'-N9	5.08	112.27	108.20
1	AA	1670	G	O4'-C1'-N9	5.08	112.27	108.20
53	B5	1560	G	O4'-C1'-N9	5.08	112.27	108.20
53	B5	2788	C	C6-N1-C1'	-5.08	114.70	120.80
1	AA	341	A	O4'-C1'-N9	5.08	112.27	108.20
1	AA	399	A	C4-C5-C6	5.08	119.54	117.00
1	AA	531	C	O4'-C1'-N1	5.08	112.27	108.20
1	AA	1134	A	O4'-C1'-N9	5.08	112.27	108.20
52	B4	131	A	C4-C5-C6	5.08	119.54	117.00
53	B5	808	A	O4'-C1'-N9	5.08	112.27	108.20
53	B5	835	G	O4'-C1'-N9	5.08	112.27	108.20
53	B5	1254	C	N3-C4-C5	-5.08	119.87	121.90
53	B5	1291	A	C5-C6-N6	-5.08	119.63	123.70
53	B5	1434	G	O4'-C1'-N9	5.08	112.27	108.20
53	B5	1561	G	C5-C6-O6	-5.08	125.55	128.60
53	B5	2521	U	O4'-C1'-N1	5.08	112.27	108.20
53	B5	2801	A	C5-C6-N1	-5.08	115.16	117.70
53	B5	2984	C	N3-C4-C5	-5.08	119.87	121.90
53	B5	2988	C	N3-C4-C5	-5.08	119.87	121.90
1	AA	586	G	C5-C6-O6	-5.08	125.55	128.60
1	AA	817	A	O4'-C1'-N9	5.08	112.27	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1081	A	C5-C6-N6	-5.08	119.64	123.70
1	AA	1185	A	C4-C5-C6	5.08	119.54	117.00
1	AA	1522	A	C5-C6-N1	-5.08	115.16	117.70
1	AA	1599	G	C5-C6-O6	-5.08	125.55	128.60
19	A7	5	A	O3'-P-O5'	5.08	113.66	104.00
53	B5	248	U	O4'-C1'-N1	5.08	112.27	108.20
53	B5	844	G	C5-C6-O6	-5.08	125.55	128.60
53	B5	904	A	C5-C6-N1	-5.08	115.16	117.70
53	B5	2458	A	C4-C5-C6	5.08	119.54	117.00
53	B5	2988	C	N3-C4-N4	5.08	121.56	118.00
53	B5	3020	U	O4'-C1'-N1	5.08	112.27	108.20
1	AA	394	C	N3-C4-N4	5.08	121.56	118.00
51	B3	65	G	O4'-C1'-N9	5.08	112.26	108.20
53	B5	761	A	C4-C5-C6	5.08	119.54	117.00
53	B5	1719	G	O4'-C1'-N9	5.08	112.26	108.20
53	B5	2952	G	P-O3'-C3'	5.08	125.80	119.70
1	AA	145	A	O4'-C1'-N9	5.08	112.26	108.20
1	AA	179	A	C4-C5-C6	5.08	119.54	117.00
1	AA	502	U	C5'-C4'-C3'	5.08	124.13	116.00
1	AA	734	A	C5-C6-N6	-5.08	119.64	123.70
1	AA	1454	C	N3-C4-N4	5.08	121.56	118.00
1	AA	1788	A	C4-C5-C6	5.08	119.54	117.00
19	A7	71	G	N9-C1'-C2'	-5.08	106.41	112.00
53	B5	473	G	O4'-C1'-N9	5.08	112.26	108.20
53	B5	495	G	C5-C6-O6	-5.08	125.55	128.60
53	B5	1003	A	C4-C5-C6	5.08	119.54	117.00
53	B5	1531	C	N3-C4-C5	-5.08	119.87	121.90
53	B5	2833	A	C4-C5-C6	5.08	119.54	117.00
1	AA	1297	A	C5-C6-N1	-5.08	115.16	117.70
19	A7	35	A	O4'-C1'-C2'	5.08	112.17	107.60
53	B5	1302	A	O4'-C1'-N9	5.08	112.26	108.20
53	B5	1852	G	O4'-C1'-N9	5.08	112.26	108.20
1	AA	351	C	N3-C4-N4	5.08	121.55	118.00
1	AA	660	G	C5-C6-O6	-5.08	125.55	128.60
1	AA	1315	G	O4'-C1'-N9	5.08	112.26	108.20
53	B5	18	G	O4'-C1'-N9	5.08	112.26	108.20
53	B5	114	A	C5-C6-N6	-5.08	119.64	123.70
53	B5	397	A	C4-C5-C6	5.08	119.54	117.00
53	B5	659	G	O4'-C1'-N9	5.08	112.26	108.20
53	B5	1745	C	N3-C4-C5	-5.08	119.87	121.90
53	B5	1854	C	N3-C4-N4	5.08	121.55	118.00
53	B5	2443	A	C5-C6-N6	-5.08	119.64	123.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	2705	A	C4-C5-C6	5.08	119.54	117.00
1	AA	218	A	O4'-C1'-N9	5.07	112.26	108.20
19	A7	3	G	C4-C5-C6	5.07	121.84	118.80
19	A7	45	G	N1-C6-O6	-5.07	116.86	119.90
53	B5	884	A	O4'-C1'-N9	5.07	112.26	108.20
53	B5	1408	G	C5-C6-O6	-5.07	125.56	128.60
53	B5	1896	A	C5-C6-N1	-5.07	115.16	117.70
53	B5	2256	A	C4-C5-C6	5.07	119.54	117.00
53	B5	2352	A	C5-C6-N1	-5.07	115.16	117.70
1	AA	594	A	O4'-C1'-N9	5.07	112.26	108.20
1	AA	1440	U	O4'-C1'-N1	5.07	112.26	108.20
1	AA	1524	A	C4-C5-C6	5.07	119.54	117.00
53	B5	349	A	C5-C6-N1	-5.07	115.16	117.70
53	B5	1055	A	O4'-C1'-N9	5.07	112.26	108.20
1	AA	1289	G	C5-C6-O6	-5.07	125.56	128.60
1	AA	1438	C	N3-C4-N4	5.07	121.55	118.00
53	B5	29	C	N3-C4-N4	5.07	121.55	118.00
53	B5	238	A	C5-C6-N6	-5.07	119.64	123.70
53	B5	329	U	O4'-C1'-N1	5.07	112.26	108.20
53	B5	1252	A	C5-C6-N1	-5.07	115.17	117.70
53	B5	1338	C	N3-C4-N4	5.07	121.55	118.00
53	B5	1588	A	C4-C5-C6	5.07	119.54	117.00
53	B5	3369	G	C4-N9-C1'	5.07	133.09	126.50
1	AA	771	A	C5-C6-N6	-5.07	119.64	123.70
1	AA	1365	G	C5-C6-O6	-5.07	125.56	128.60
53	B5	1248	C	N3-C4-N4	5.07	121.55	118.00
53	B5	1893	A	C4-C5-C6	5.07	119.53	117.00
53	B5	2879	C	N3-C4-N4	5.07	121.55	118.00
53	B5	3247	G	C5-C6-O6	-5.07	125.56	128.60
1	AA	350	U	O4'-C1'-N1	5.07	112.25	108.20
1	AA	365	G	O4'-C1'-N9	5.07	112.25	108.20
1	AA	445	A	C4-C5-C6	5.07	119.53	117.00
1	AA	903	G	O4'-C1'-N9	5.07	112.25	108.20
1	AA	954	A	C5-C6-N6	-5.07	119.65	123.70
1	AA	1150	G	C5-C6-O6	-5.07	125.56	128.60
1	AA	1319	A	O4'-C1'-N9	5.07	112.25	108.20
1	AA	1516	C	P-O5'-C5'	-5.07	112.79	120.90
1	AA	1612	A	C5-C6-N1	-5.07	115.17	117.70
52	B4	1	A	C4-C5-C6	5.07	119.53	117.00
53	B5	318	A	C4-C5-C6	5.07	119.53	117.00
53	B5	363	G	O4'-C1'-N9	5.07	112.25	108.20
53	B5	653	A	C4-C5-C6	5.07	119.53	117.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	1000	C	N3-C4-N4	5.07	121.55	118.00
53	B5	2658	G	O4'-C1'-N9	5.07	112.25	108.20
53	B5	2674	A	C5-C6-N6	-5.07	119.65	123.70
53	B5	2732	G	O4'-C1'-N9	5.07	112.25	108.20
53	B5	2857	C	N3-C4-N4	5.07	121.55	118.00
53	B5	2991	A	C5-C6-N1	-5.07	115.17	117.70
53	B5	3245	A	O4'-C1'-N9	5.07	112.25	108.20
1	AA	948	C	N3-C4-N4	5.07	121.55	118.00
1	AA	962	A	C4-C5-C6	5.07	119.53	117.00
1	AA	1148	A	O4'-C1'-N9	5.07	112.25	108.20
1	AA	1237	G	C5-C6-O6	-5.07	125.56	128.60
52	B4	40	A	C4-C5-C6	5.07	119.53	117.00
53	B5	304	G	O4'-C1'-N9	5.07	112.25	108.20
53	B5	711	A	C4-C5-C6	5.07	119.53	117.00
53	B5	1482	A	C4-C5-C6	5.07	119.53	117.00
53	B5	1490	A	O4'-C1'-N9	5.07	112.25	108.20
53	B5	1671	C	N3-C4-N4	5.07	121.55	118.00
53	B5	2656	A	C5-C6-N1	-5.07	115.17	117.70
53	B5	3308	C	N3-C4-N4	5.07	121.55	118.00
1	AA	962	A	C5-C6-N6	-5.06	119.65	123.70
1	AA	1451	G	C5-C6-O6	-5.06	125.56	128.60
19	A7	51	G	N9-C4-C5	5.06	107.43	105.40
53	B5	135	C	C6-N1-C2	-5.06	118.28	120.30
53	B5	728	G	O4'-C1'-N9	5.06	112.25	108.20
53	B5	1497	C	N3-C4-C5	-5.06	119.87	121.90
1	AA	199	G	C5-C6-O6	-5.06	125.56	128.60
1	AA	724	C	N3-C4-C5	-5.06	119.88	121.90
1	AA	1140	A	C4-C5-C6	5.06	119.53	117.00
1	AA	1310	A	C5-C6-N6	-5.06	119.65	123.70
53	B5	1673	G	C5-C6-O6	-5.06	125.56	128.60
53	B5	2267	C	N3-C4-C5	-5.06	119.88	121.90
53	B5	3279	A	C4-C5-C6	5.06	119.53	117.00
53	B5	3347	A	C4-C5-C6	5.06	119.53	117.00
1	AA	858	G	N9-C1'-C2'	5.06	120.58	114.00
1	AA	1649	A	O4'-C1'-N9	5.06	112.25	108.20
53	B5	2443	A	O4'-C1'-N9	5.06	112.25	108.20
1	AA	540	G	C5-C6-O6	-5.06	125.56	128.60
1	AA	731	C	N3-C4-N4	5.06	121.54	118.00
1	AA	1744	A	C4-C5-C6	5.06	119.53	117.00
52	B4	133	G	C5-C6-O6	-5.06	125.56	128.60
53	B5	268	A	C4-C5-C6	5.06	119.53	117.00
53	B5	781	G	O4'-C1'-N9	5.06	112.25	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	2208	A	C5-C6-N1	-5.06	115.17	117.70
53	B5	2400	G	O4'-C1'-N9	5.06	112.25	108.20
53	B5	2646	C	N3-C4-C5	-5.06	119.88	121.90
53	B5	3179	U	P-O3'-C3'	5.06	125.77	119.70
1	AA	552	G	O4'-C1'-N9	5.06	112.25	108.20
1	AA	696	C	N3-C4-C5	-5.06	119.88	121.90
1	AA	1085	A	C4-C5-C6	5.06	119.53	117.00
52	B4	70	A	C4-C5-C6	5.06	119.53	117.00
52	B4	94	C	N3-C4-C5	-5.06	119.88	121.90
53	B5	61	A	C4-C5-C6	5.06	119.53	117.00
53	B5	101	G	C5-C6-O6	-5.06	125.57	128.60
53	B5	709	A	C5-C6-N6	-5.06	119.66	123.70
53	B5	1295	G	C5-C6-O6	-5.06	125.57	128.60
53	B5	1733	G	O4'-C1'-N9	5.06	112.25	108.20
53	B5	2178	A	C5-C6-N1	-5.06	115.17	117.70
53	B5	2469	G	O4'-C1'-N9	5.06	112.25	108.20
53	B5	3048	A	C5-C6-N1	-5.06	115.17	117.70
1	AA	518	A	O4'-C1'-N9	5.06	112.25	108.20
1	AA	1639	C	N3-C4-N4	5.06	121.54	118.00
4	AD	78	ARG	NE-CZ-NH2	-5.06	117.77	120.30
19	A7	52	U	C6-N1-C2	-5.06	117.97	121.00
52	B4	130	C	N3-C4-N4	5.06	121.54	118.00
53	B5	2645	G	O4'-C1'-N9	5.06	112.25	108.20
53	B5	2911	A	C4-C5-C6	5.06	119.53	117.00
53	B5	3112	G	O4'-C1'-N9	5.06	112.25	108.20
1	AA	188	A	C4-C5-C6	5.05	119.53	117.00
1	AA	1166	G	O4'-C1'-N9	5.05	112.24	108.20
1	AA	1682	U	O4'-C1'-N1	5.05	112.24	108.20
52	B4	54	A	C4-C5-C6	5.05	119.53	117.00
53	B5	701	G	O4'-C1'-N9	5.05	112.24	108.20
53	B5	1120	A	C4-C5-C6	5.05	119.53	117.00
53	B5	3374	U	O4'-C1'-N1	5.05	112.24	108.20
1	AA	338	C	N3-C4-C5	-5.05	119.88	121.90
51	B3	27	A	O4'-C1'-N9	5.05	112.24	108.20
52	B4	149	A	C4-C5-C6	5.05	119.53	117.00
53	B5	144	A	C4-C5-C6	5.05	119.53	117.00
53	B5	1193	A	O4'-C1'-N9	5.05	112.24	108.20
53	B5	1608	C	N3-C4-N4	5.05	121.54	118.00
53	B5	1857	C	N3-C4-C5	-5.05	119.88	121.90
53	B5	2901	G	C5-C6-O6	-5.05	125.57	128.60
53	B5	2926	A	C5-C6-N1	-5.05	115.17	117.70
1	AA	226	A	O4'-C1'-N9	5.05	112.24	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1334	A	O4'-C1'-N9	5.05	112.24	108.20
1	AA	1572	G	C5-C6-O6	-5.05	125.57	128.60
53	B5	445	G	C5-C6-O6	-5.05	125.57	128.60
53	B5	978	C	P-O3'-C3'	5.05	125.76	119.70
53	B5	1086	C	N3-C4-N4	5.05	121.54	118.00
1	AA	295	A	O4'-C1'-N9	5.05	112.24	108.20
1	AA	605	A	O4'-C1'-N9	5.05	112.24	108.20
1	AA	859	A	C4'-C3'-C2'	5.05	107.65	102.60
1	AA	1224	G	O4'-C1'-N9	5.05	112.24	108.20
1	AA	1789	A	C5-C6-N6	-5.05	119.66	123.70
53	B5	973	A	C4-C5-C6	5.05	119.53	117.00
53	B5	1118	C	N3-C4-N4	5.05	121.53	118.00
53	B5	1487	G	O4'-C1'-N9	5.05	112.24	108.20
53	B5	1525	G	O4'-C1'-N9	5.05	112.24	108.20
53	B5	1700	G	C5-C6-O6	-5.05	125.57	128.60
53	B5	3017	A	C4-C5-C6	5.05	119.53	117.00
53	B5	3169	U	O4'-C1'-N1	5.05	112.24	108.20
1	AA	188	A	C5-C6-N6	-5.05	119.66	123.70
53	B5	2225	U	O4'-C1'-N1	5.05	112.24	108.20
53	B5	2468	A	C5-C6-N6	-5.05	119.66	123.70
1	AA	89	G	O4'-C1'-N9	5.05	112.24	108.20
1	AA	274	G	C5-C6-O6	-5.05	125.57	128.60
1	AA	1201	C	N3-C4-N4	5.05	121.53	118.00
1	AA	1728	A	O4'-C1'-N9	5.05	112.24	108.20
19	A7	20	G	N9-C4-C5	5.05	107.42	105.40
53	B5	265	A	C4-C5-C6	5.05	119.52	117.00
53	B5	304	G	C5-C6-O6	-5.05	125.57	128.60
53	B5	593	A	C5-C6-N1	-5.05	115.18	117.70
53	B5	739	G	C5-C6-O6	-5.05	125.57	128.60
53	B5	743	C	N3-C4-N4	5.05	121.53	118.00
53	B5	1236	G	N3-C2-N2	5.05	123.43	119.90
53	B5	1339	C	N3-C4-N4	5.05	121.53	118.00
53	B5	1704	A	O4'-C1'-N9	5.05	112.24	108.20
53	B5	1741	A	C5-C6-N6	-5.05	119.66	123.70
53	B5	1750	A	C5-C6-N6	-5.05	119.66	123.70
53	B5	2699	G	C5-C6-O6	-5.05	125.57	128.60
1	AA	755	A	C5-C6-N6	-5.04	119.66	123.70
53	B5	12	A	O4'-C1'-N9	5.04	112.24	108.20
53	B5	1238	C	N3-C4-N4	5.04	121.53	118.00
1	AA	510	G	C5-C6-O6	-5.04	125.57	128.60
1	AA	914	A	C5-C6-N6	-5.04	119.67	123.70
1	AA	1552	U	P-O5'-C5'	-5.04	112.83	120.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
19	A7	15	G	C8-N9-C1'	5.04	133.56	127.00
52	B4	131	A	C5-C6-N1	-5.04	115.18	117.70
53	B5	808	A	C4-C5-C6	5.04	119.52	117.00
53	B5	3165	A	O4'-C1'-N9	5.04	112.23	108.20
1	AA	1	U	C6-N1-C1'	-5.04	114.14	121.20
1	AA	939	A	C5-C6-N1	-5.04	115.18	117.70
1	AA	1170	C	N3-C4-C5	-5.04	119.88	121.90
52	B4	45	C	N3-C4-C5	-5.04	119.88	121.90
53	B5	208	C	N3-C4-N4	5.04	121.53	118.00
53	B5	352	A	C4-C5-C6	5.04	119.52	117.00
53	B5	661	G	O4'-C1'-N9	5.04	112.23	108.20
53	B5	887	G	C5-C6-O6	-5.04	125.58	128.60
53	B5	2243	A	C5-C6-N6	-5.04	119.67	123.70
53	B5	2622	C	N3-C4-C5	-5.04	119.88	121.90
53	B5	2721	A	C4-C5-C6	5.04	119.52	117.00
53	B5	2989	U	O4'-C1'-N1	5.04	112.23	108.20
53	B5	3303	G	C5-C6-O6	-5.04	125.58	128.60
53	B5	533	A	O4'-C1'-N9	5.04	112.23	108.20
53	B5	1779	C	N3-C4-C5	-5.04	119.88	121.90
53	B5	2598	G	C5-C6-O6	-5.04	125.58	128.60
1	AA	67	A	C4-C5-C6	5.04	119.52	117.00
51	B3	92	A	C5-C6-N6	-5.04	119.67	123.70
53	B5	347	G	C5-C6-O6	-5.04	125.58	128.60
53	B5	715	A	C1'-O4'-C4'	-5.04	105.87	109.90
53	B5	1797	A	C4-C5-C6	5.04	119.52	117.00
1	AA	954	A	O4'-C1'-N9	5.04	112.23	108.20
1	AA	1281	C	N3-C4-C5	-5.04	119.89	121.90
1	AA	1456	G	C5-C6-O6	-5.04	125.58	128.60
1	AA	1468	C	N3-C4-N4	5.04	121.53	118.00
1	AA	1725	G	C5-C6-O6	-5.04	125.58	128.60
53	B5	17	G	C5-C6-O6	-5.04	125.58	128.60
53	B5	878	G	C5-C6-O6	-5.04	125.58	128.60
53	B5	1797	A	O4'-C1'-N9	5.04	112.23	108.20
53	B5	2196	C	N3-C4-C5	-5.04	119.89	121.90
1	AA	509	G	C5-C6-O6	-5.04	125.58	128.60
1	AA	864	U	OP1-P-OP2	-5.04	112.05	119.60
1	AA	1245	U	O4'-C1'-N1	5.04	112.23	108.20
1	AA	1635	C	O3'-P-O5'	5.04	113.57	104.00
53	B5	255	A	C4-C5-C6	5.04	119.52	117.00
53	B5	690	A	C4-C5-C6	5.04	119.52	117.00
53	B5	1239	C	N3-C4-N4	5.04	121.53	118.00
53	B5	2144	A	C4-C5-C6	5.04	119.52	117.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	2432	A	O4'-C1'-N9	5.04	112.23	108.20
53	B5	2491	A	C4-C5-C6	5.04	119.52	117.00
53	B5	2534	G	C5-C6-O6	-5.04	125.58	128.60
53	B5	2987	A	C4-C5-C6	5.04	119.52	117.00
53	B5	3254	G	C5-C6-O6	-5.04	125.58	128.60
1	AA	22	A	O4'-C1'-N9	5.03	112.23	108.20
1	AA	202	A	O4'-C1'-N9	5.03	112.23	108.20
1	AA	531	C	N3-C4-N4	5.03	121.52	118.00
19	A7	60	C	C4-C5-C6	-5.03	114.88	117.40
53	B5	229	G	O4'-C1'-N9	5.03	112.23	108.20
53	B5	975	G	O4'-C1'-N9	5.03	112.23	108.20
53	B5	1321	G	C5-C6-O6	-5.03	125.58	128.60
53	B5	2114	C	N3-C4-N4	5.03	121.52	118.00
53	B5	2479	C	N3-C4-C5	-5.03	119.89	121.90
1	AA	1279	U	O4'-C1'-N1	5.03	112.23	108.20
19	A7	14	A	O4'-C1'-N9	5.03	112.23	108.20
33	BI	159	PHE	CB-CG-CD2	-5.03	117.28	120.80
53	B5	397	A	C5-C6-N6	-5.03	119.67	123.70
53	B5	817	A	C4-C5-C6	5.03	119.52	117.00
53	B5	1225	A	O4'-C1'-N9	5.03	112.22	108.20
53	B5	3323	A	C5-C6-N6	-5.03	119.67	123.70
1	AA	202	A	C4-C5-C6	5.03	119.52	117.00
1	AA	783	G	C5-C6-O6	-5.03	125.58	128.60
1	AA	1099	G	O4'-C1'-N9	5.03	112.22	108.20
1	AA	1580	U	O5'-P-OP2	-5.03	101.17	105.70
53	B5	807	A	O4'-C1'-N9	5.03	112.22	108.20
53	B5	842	G	O4'-C1'-N9	5.03	112.22	108.20
53	B5	1159	A	C5-C6-N6	-5.03	119.68	123.70
53	B5	1184	A	C5-C6-N1	-5.03	115.19	117.70
53	B5	1933	A	C4-C5-C6	5.03	119.52	117.00
53	B5	2798	C	N3-C4-N4	5.03	121.52	118.00
53	B5	2824	G	O4'-C1'-N9	5.03	112.22	108.20
53	B5	3210	U	P-O3'-C3'	5.03	125.74	119.70
1	AA	1088	A	C4-C5-C6	5.03	119.51	117.00
1	AA	1397	A	O4'-C1'-N9	5.03	112.22	108.20
53	B5	1133	A	C5-C6-N6	-5.03	119.68	123.70
53	B5	1496	C	N3-C4-N4	5.03	121.52	118.00
53	B5	2105	G	C5-C6-O6	-5.03	125.58	128.60
1	AA	1328	A	C5-C6-N6	-5.03	119.68	123.70
1	AA	1631	A	C4'-C3'-C2'	5.03	107.63	102.60
1	AA	1690	G	C5-C6-O6	-5.03	125.58	128.60
18	AT	78	LYS	N-CA-C	5.03	124.57	111.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	578	A	C4-C5-C6	5.03	119.51	117.00
53	B5	1043	C	N3-C4-C5	-5.03	119.89	121.90
53	B5	1423	C	N3-C4-C5	-5.03	119.89	121.90
53	B5	2259	A	O4'-C1'-N9	5.03	112.22	108.20
53	B5	2356	A	O4'-C1'-N9	5.03	112.22	108.20
53	B5	3164	C	O4'-C1'-N1	5.03	112.22	108.20
1	AA	470	A	C4-C5-C6	5.03	119.51	117.00
1	AA	1485	A	C4-C5-C6	5.03	119.51	117.00
51	B3	71	C	N3-C4-N4	5.03	121.52	118.00
53	B5	41	G	C5-C6-O6	-5.03	125.58	128.60
53	B5	65	A	C5-C6-N6	-5.03	119.68	123.70
53	B5	578	A	C5-C6-N1	-5.03	115.19	117.70
53	B5	827	A	C4-C5-C6	5.03	119.51	117.00
53	B5	973	A	O4'-C1'-N9	5.03	112.22	108.20
53	B5	1221	A	C5-C6-N6	-5.03	119.68	123.70
53	B5	2156	C	N3-C4-N4	5.03	121.52	118.00
1	AA	928	A	O4'-C1'-N9	5.02	112.22	108.20
53	B5	1569	U	O4'-C1'-N1	5.02	112.22	108.20
1	AA	91	G	C5-C6-O6	-5.02	125.59	128.60
1	AA	1215	A	C5-C6-N6	-5.02	119.68	123.70
1	AA	1294	G	C5-C6-O6	-5.02	125.59	128.60
1	AA	1492	C	N3-C4-N4	5.02	121.52	118.00
53	B5	41	G	O4'-C1'-N9	5.02	112.22	108.20
53	B5	307	A	C5-C6-N1	-5.02	115.19	117.70
53	B5	1696	A	C5-C6-N6	-5.02	119.68	123.70
53	B5	2166	A	O4'-C1'-N9	5.02	112.22	108.20
53	B5	3184	U	O4'-C1'-N1	5.02	112.22	108.20
53	B5	238	A	O4'-C1'-N9	5.02	112.22	108.20
53	B5	1867	A	C4-C5-C6	5.02	119.51	117.00
53	B5	2220	A	C4-C5-C6	5.02	119.51	117.00
53	B5	2228	A	C4-C5-C6	5.02	119.51	117.00
1	AA	1772	G	O4'-C1'-N9	5.02	112.22	108.20
27	BC	67	PHE	CB-CG-CD2	5.02	124.31	120.80
51	B3	42	A	C4-C5-C6	5.02	119.51	117.00
51	B3	102	C	N3-C4-C5	-5.02	119.89	121.90
53	B5	278	U	O4'-C1'-N1	5.02	112.22	108.20
53	B5	625	G	C5-C6-O6	-5.02	125.59	128.60
53	B5	973	A	C5-C6-N1	-5.02	115.19	117.70
53	B5	2234	G	C5-C6-O6	-5.02	125.59	128.60
53	B5	2350	C	N3-C4-N4	5.02	121.51	118.00
53	B5	2708	C	N3-C4-C5	-5.02	119.89	121.90
53	B5	2847	A	O4'-C1'-N9	5.02	112.22	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	3018	C	N3-C4-C5	-5.02	119.89	121.90
53	B5	3032	A	C5-C6-N6	-5.02	119.68	123.70
53	B5	3033	A	C4-C5-C6	5.02	119.51	117.00
1	AA	965	A	C5-C6-N6	-5.02	119.69	123.70
1	AA	1122	A	C4-C5-C6	5.02	119.51	117.00
1	AA	1795	A	C4-C5-C6	5.02	119.51	117.00
19	A7	15	G	O4'-C4'-C3'	5.02	110.11	106.10
53	B5	26	A	C4-C5-C6	5.02	119.51	117.00
53	B5	717	C	N3-C4-N4	5.02	121.51	118.00
53	B5	1225	A	C5-C6-N6	-5.02	119.69	123.70
53	B5	1304	A	O4'-C1'-N9	5.02	112.21	108.20
53	B5	2874	G	C5-C6-O6	-5.02	125.59	128.60
1	AA	190	C	N3-C4-N4	5.02	121.51	118.00
53	B5	745	C	N3-C4-C5	-5.02	119.89	121.90
53	B5	1205	A	O4'-C1'-N9	5.02	112.21	108.20
53	B5	1418	A	C5-C6-N6	-5.02	119.69	123.70
53	B5	1834	U	O4'-C1'-N1	5.02	112.21	108.20
53	B5	2932	U	O4'-C1'-N1	5.02	112.21	108.20
53	B5	3073	A	C4-C5-C6	5.02	119.51	117.00
1	AA	570	A	C4-C5-C6	5.01	119.51	117.00
1	AA	1503	A	C5-C6-N6	-5.01	119.69	123.70
4	AD	78	ARG	NE-CZ-NH1	5.01	122.81	120.30
51	B3	101	A	C4-C5-C6	5.01	119.51	117.00
53	B5	408	A	O4'-C1'-N9	5.01	112.21	108.20
53	B5	455	C	P-O3'-C3'	5.01	125.72	119.70
53	B5	735	A	C5-C6-N6	-5.01	119.69	123.70
53	B5	1804	A	C5-C6-N6	-5.01	119.69	123.70
53	B5	2736	A	O4'-C1'-N9	5.01	112.21	108.20
53	B5	2977	G	C5-C6-O6	-5.01	125.59	128.60
53	B5	3181	G	C5-C6-O6	-5.01	125.59	128.60
53	B5	3370	A	C4-C5-C6	5.01	119.51	117.00
1	AA	1315	G	C5-C6-O6	-5.01	125.59	128.60
1	AA	1540	G	C5-C6-O6	-5.01	125.59	128.60
1	AA	1770	C	N3-C4-N4	5.01	121.51	118.00
53	B5	963	G	O4'-C1'-N9	5.01	112.21	108.20
53	B5	1424	C	N3-C4-C5	-5.01	119.89	121.90
53	B5	2580	A	C5-C6-N6	-5.01	119.69	123.70
53	B5	2971	A	O4'-C1'-N9	5.01	112.21	108.20
53	B5	3168	A	C4-C5-C6	5.01	119.51	117.00
1	AA	6	G	O4'-C1'-N9	5.01	112.21	108.20
1	AA	301	A	C5-C6-N6	-5.01	119.69	123.70
1	AA	874	C	N3-C4-N4	5.01	121.51	118.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1038	A	C5-C6-N6	-5.01	119.69	123.70
51	B3	70	A	C4-C5-C6	5.01	119.50	117.00
53	B5	846	A	C5-C6-N6	-5.01	119.69	123.70
53	B5	923	C	N3-C4-N4	5.01	121.51	118.00
53	B5	1153	A	C5-C6-N6	-5.01	119.69	123.70
53	B5	1507	G	N1-C6-O6	5.01	122.91	119.90
53	B5	1835	A	O4'-C1'-N9	5.01	112.21	108.20
53	B5	2306	C	O4'-C1'-N1	5.01	112.21	108.20
53	B5	2649	A	C4-C5-C6	5.01	119.51	117.00
53	B5	3026	G	O4'-C1'-N9	5.01	112.21	108.20
53	B5	3135	U	P-O3'-C3'	5.01	125.71	119.70
1	AA	334	G	C5-C6-O6	-5.01	125.59	128.60
1	AA	696	C	N3-C4-N4	5.01	121.51	118.00
1	AA	1164	G	C5-C6-O6	-5.01	125.59	128.60
1	AA	1775	G	O4'-C1'-N9	5.01	112.21	108.20
1	AA	1791	G	C5-C6-O6	-5.01	125.59	128.60
19	A7	44	A	C6-C5-N7	5.01	135.81	132.30
53	B5	580	C	N3-C4-C5	-5.01	119.90	121.90
53	B5	1009	A	O4'-C1'-N9	5.01	112.21	108.20
53	B5	1290	A	O4'-C1'-N9	5.01	112.21	108.20
53	B5	1806	A	C4-C5-C6	5.01	119.50	117.00
53	B5	2703	A	C4-C5-C6	5.01	119.50	117.00
53	B5	3002	C	N3-C4-C5	-5.01	119.90	121.90
53	B5	3347	A	C5-C6-N6	-5.01	119.69	123.70
1	AA	1330	C	N3-C4-C5	-5.01	119.90	121.90
53	B5	2296	A	C5-C6-N1	-5.01	115.20	117.70
53	B5	3207	C	N3-C4-N4	5.01	121.50	118.00
53	B5	379	C	N3-C4-C5	-5.01	119.90	121.90
53	B5	1947	G	C4-N9-C1'	5.01	133.01	126.50
53	B5	2312	A	C4-C5-C6	5.01	119.50	117.00
53	B5	2779	A	C4-C5-C6	5.01	119.50	117.00
53	B5	2952	G	O4'-C1'-N9	5.01	112.20	108.20
53	B5	3016	A	P-O3'-C3'	5.01	125.71	119.70
53	B5	3133	C	N3-C4-C5	-5.01	119.90	121.90
53	B5	13	A	C4-C5-C6	5.00	119.50	117.00
53	B5	466	G	O4'-C1'-N9	5.00	112.20	108.20
53	B5	598	A	O4'-C1'-N9	5.00	112.20	108.20
53	B5	2106	A	C5-C6-N1	-5.00	115.20	117.70
53	B5	3218	A	C4-C5-C6	5.00	119.50	117.00
1	AA	341	A	C4-C5-C6	5.00	119.50	117.00
1	AA	1326	A	C5-C6-N6	-5.00	119.70	123.70
53	B5	408	A	C5-C6-N6	-5.00	119.70	123.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	B5	2157	G	C5-C6-O6	-5.00	125.60	128.60
1	AA	1160	A	C4-C5-C6	5.00	119.50	117.00
51	B3	2	G	C5-C6-O6	-5.00	125.60	128.60
51	B3	102	C	N3-C4-N4	5.00	121.50	118.00
53	B5	747	A	O4'-C1'-N9	5.00	112.20	108.20
53	B5	2251	G	P-O3'-C3'	5.00	125.70	119.70
53	B5	2447	A	C4-C5-C6	5.00	119.50	117.00
53	B5	3263	G	C5-C6-O6	-5.00	125.60	128.60
53	B5	3279	A	O4'-C1'-N9	5.00	112.20	108.20

All (45) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	AA	109	G	C3'
1	AA	933	C	C2'
7	AH	34	ILE	CB
26	BB	161	ASP	CA
29	BE	216	GLU	CA
30	BF	83	LEU	CA
31	BG	141	ALA	CA
31	BG	142	LEU	CA
36	BL	56	LYS	CA
48	BX	70	ALA	CA
48	BX	80	ALA	CA
51	B3	70	A	C4'
53	B5	114	A	C3'
53	B5	279	U	C3'
53	B5	355	A	C4'
53	B5	477	A	C3'
53	B5	481	U	C3'
53	B5	554	A	C3'
53	B5	596	U	C3'
53	B5	705	A	C3'
53	B5	711	A	C2'
53	B5	735	A	C2',C3'
53	B5	786	A	C3'
53	B5	970	A	C3'
53	B5	1347	U	C3'
53	B5	1583	A	C2',C3'
53	B5	1590	G	C3'
53	B5	1644	C	C2'
53	B5	1737	U	C3'

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Mol	Chain	Res	Type	Atom
53	B5	1947	G	C1'
53	B5	2214	A	C3'
53	B5	2224	A	C3'
53	B5	2225	U	C3'
53	B5	2229	A	C3'
53	B5	2469	G	C2',C1'
53	B5	2477	G	C1'
53	B5	2999	U	C2'
53	B5	3051	U	C4'
53	B5	3142	A	C3'
53	B5	3246	G	C4'
53	B5	3303	G	C2'
53	B5	3304	U	C3'

All (285) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
19	A7	1	G	Sidechain
19	A7	11	C	Sidechain
19	A7	14	A	Sidechain
19	A7	15	G	Sidechain
19	A7	18	G	Sidechain
19	A7	19	G	Sidechain
19	A7	2	C	Sidechain
19	A7	20	G	Sidechain
19	A7	21	A	Sidechain
19	A7	23	A	Sidechain
19	A7	24	G	Sidechain
19	A7	27	C	Sidechain
19	A7	29	A	Sidechain
19	A7	3	G	Sidechain
19	A7	30	G	Sidechain
19	A7	31	A	Sidechain
19	A7	36	A	Sidechain
19	A7	4	G	Sidechain
19	A7	41	U	Sidechain
19	A7	42	G	Sidechain
19	A7	43	G	Sidechain
19	A7	45	G	Sidechain
19	A7	47	U	Sidechain
19	A7	5	A	Sidechain
19	A7	50	U	Sidechain

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Mol	Chain	Res	Type	Group
19	A7	51	G	Sidechain
19	A7	52	U	Sidechain
19	A7	53	G	Sidechain
19	A7	56	C	Sidechain
19	A7	57	G	Sidechain
19	A7	59	U	Sidechain
19	A7	61	C	Sidechain
19	A7	63	C	Sidechain
19	A7	65	G	Sidechain
19	A7	66	A	Sidechain
19	A7	7	U	Sidechain
19	A7	70	C	Sidechain
19	A7	71	G	Sidechain
19	A7	72	C	Sidechain
19	A7	73	A	Sidechain
19	A7	74	C	Sidechain
19	A7	75	C	Sidechain
19	A7	76	A	Sidechain
19	A7	9	A	Sidechain
1	AA	1089	A	Sidechain
1	AA	1196	G	Sidechain
1	AA	1200	A	Sidechain
1	AA	1202	U	Sidechain
1	AA	1237	G	Sidechain
1	AA	1303	C	Sidechain
1	AA	1309	A	Sidechain
1	AA	1327	G	Sidechain
1	AA	1440	U	Sidechain
1	AA	1502	G	Sidechain
1	AA	1514	A	Sidechain
1	AA	1520	U	Sidechain
1	AA	1551	G	Sidechain
1	AA	1583	U	Sidechain
1	AA	1584	A	Sidechain
1	AA	1586	G	Sidechain
1	AA	1766	G	Sidechain
1	AA	225	A	Sidechain
1	AA	244	A	Sidechain
1	AA	474	A	Sidechain
1	AA	475	A	Sidechain
1	AA	476	U	Sidechain
1	AA	529	A	Sidechain

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Mol	Chain	Res	Type	Group
1	AA	80	A	Sidechain
1	AA	843	U	Sidechain
1	AA	855	A	Sidechain
1	AA	856	A	Sidechain
1	AA	877	G	Sidechain
2	AB	84	ARG	Sidechain
4	AD	149	ARG	Sidechain
4	AD	78	ARG	Sidechain
4	AD	79	ARG	Sidechain
5	AE	226	THR	Peptide
8	AI	135	ARG	Sidechain
8	AI	142	TYR	Sidechain
8	AI	96	TYR	Sidechain
9	AJ	23	ARG	Sidechain
9	AJ	53	LYS	Peptide
9	AJ	90	TYR	Sidechain
10	AK	132	ARG	Sidechain,Peptide
10	AK	48	VAL	Peptide
10	AK	52	ARG	Sidechain
12	AM	120	ARG	Peptide
12	AM	137	HIS	Peptide
12	AM	143	ARG	Sidechain
13	AN	13	ARG	Sidechain
13	AN	14	PHE	Sidechain
14	AO	114	ARG	Sidechain
14	AO	121	ARG	Sidechain
14	AO	124	ARG	Sidechain
14	AO	135	LEU	Peptide
20	B0	45	ARG	Sidechain
52	B4	132	G	Sidechain
52	B4	147	U	Sidechain
52	B4	41	A	Sidechain
52	B4	44	A	Sidechain
52	B4	62	C	Sidechain
52	B4	64	U	Sidechain
53	B5	1	G	Sidechain
53	B5	10	C	Sidechain
53	B5	1005	G	Sidechain
53	B5	1026	A	Sidechain
53	B5	1104	G	Sidechain
53	B5	1132	C	Sidechain
53	B5	1169	A	Sidechain

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Mol	Chain	Res	Type	Group
53	B5	118	U	Sidechain
53	B5	1181	U	Sidechain
53	B5	1200	A	Sidechain
53	B5	1236	G	Sidechain
53	B5	1239	C	Sidechain
53	B5	1240	A	Sidechain
53	B5	1257	C	Sidechain
53	B5	1290	A	Sidechain
53	B5	130	A	Sidechain
53	B5	1346	G	Sidechain
53	B5	1347	U	Sidechain
53	B5	135	C	Sidechain
53	B5	1357	G	Sidechain
53	B5	1362	G	Sidechain
53	B5	1368	U	Sidechain
53	B5	1378	U	Sidechain
53	B5	1385	C	Sidechain
53	B5	1417	G	Sidechain
53	B5	1421	G	Sidechain
53	B5	1496	C	Sidechain
53	B5	1510	G	Sidechain
53	B5	1544	G	Sidechain
53	B5	1558	A	Sidechain
53	B5	1622	U	Sidechain
53	B5	1658	G	Sidechain
53	B5	17	G	Sidechain
53	B5	1714	A	Sidechain
53	B5	1734	G	Sidechain
53	B5	1783	U	Sidechain
53	B5	1790	G	Sidechain
53	B5	18	G	Sidechain
53	B5	1811	G	Sidechain
53	B5	1812	G	Sidechain
53	B5	1831	U	Sidechain
53	B5	1874	A	Sidechain
53	B5	1896	A	Sidechain
53	B5	1927	G	Sidechain
53	B5	2112	U	Sidechain
53	B5	2130	G	Sidechain
53	B5	2133	U	Sidechain
53	B5	216	G	Sidechain
53	B5	2167	A	Sidechain

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Mol	Chain	Res	Type	Group
53	B5	2176	U	Sidechain
53	B5	2178	A	Sidechain
53	B5	2180	G	Sidechain
53	B5	2185	G	Sidechain
53	B5	2186	U	Sidechain
53	B5	2224	A	Sidechain
53	B5	2300	G	Sidechain
53	B5	2356	A	Sidechain
53	B5	2394	G	Sidechain
53	B5	2396	G	Sidechain
53	B5	2448	G	Sidechain
53	B5	2468	A	Sidechain
53	B5	2481	G	Sidechain
53	B5	2482	U	Sidechain
53	B5	251	G	Sidechain
53	B5	2548	C	Sidechain
53	B5	2586	G	Sidechain
53	B5	2593	A	Sidechain
53	B5	2607	G	Sidechain
53	B5	2621	G	Sidechain
53	B5	2677	G	Sidechain
53	B5	2681	U	Sidechain
53	B5	2696	A	Sidechain
53	B5	270	U	Sidechain
53	B5	2744	U	Sidechain
53	B5	2751	G	Sidechain
53	B5	2768	U	Sidechain
53	B5	2774	C	Sidechain
53	B5	2789	U	Sidechain
53	B5	2839	G	Sidechain
53	B5	2876	C	Sidechain
53	B5	2901	G	Sidechain
53	B5	2943	G	Sidechain
53	B5	2951	G	Sidechain
53	B5	3003	G	Sidechain
53	B5	3009	G	Sidechain
53	B5	3032	A	Sidechain
53	B5	3049	A	Sidechain
53	B5	3144	G	Sidechain
53	B5	3172	G	Sidechain
53	B5	3303	G	Sidechain
53	B5	3337	G	Sidechain

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Mol	Chain	Res	Type	Group
53	B5	337	G	Sidechain
53	B5	366	A	Sidechain
53	B5	40	A	Sidechain
53	B5	494	G	Sidechain
53	B5	495	G	Sidechain
53	B5	496	C	Sidechain
53	B5	498	A	Sidechain
53	B5	584	G	Sidechain
53	B5	619	A	Sidechain
53	B5	652	G	Sidechain
53	B5	662	U	Sidechain
53	B5	709	A	Sidechain
53	B5	8	C	Sidechain
53	B5	847	A	Sidechain
53	B5	855	U	Sidechain
53	B5	859	G	Sidechain
53	B5	860	G	Sidechain
53	B5	895	A	Sidechain
53	B5	975	G	Sidechain
53	B5	977	U	Sidechain
53	B5	992	G	Sidechain
23	B8	115	ALA	Peptide
23	B8	5	ARG	Sidechain
23	B8	61	ARG	Sidechain
24	B9	18	TYR	Sidechain
24	B9	34	HIS	Peptide
24	B9	43	GLY	Peptide
26	BB	150	LEU	Peptide
26	BB	161	ASP	Peptide
26	BB	163	ARG	Sidechain
26	BB	242	ARG	Sidechain
27	BC	117	ARG	Sidechain
28	BD	202	ARG	Sidechain
28	BD	31	ARG	Sidechain
28	BD	98	ARG	Sidechain
29	BE	112	ARG	Sidechain
29	BE	192	PRO	Peptide
29	BE	198	TYR	Sidechain
29	BE	213	ASP	Peptide
29	BE	219	PHE	Peptide
29	BE	85	ARG	Sidechain
29	BE	86	TYR	Sidechain

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Mol	Chain	Res	Type	Group
31	BG	142	LEU	Peptide
31	BG	173	MET	Peptide
32	BH	173	ARG	Sidechain
32	BH	23	ARG	Sidechain
32	BH	69	ARG	Sidechain
33	BI	69	ARG	Sidechain
33	BI	7	ARG	Sidechain
34	BJ	47	GLN	Peptide
34	BJ	61	ARG	Sidechain
34	BJ	92	ARG	Sidechain
35	BK	16	ARG	Sidechain
35	BK	90	ARG	Sidechain
36	BL	153	ASP	Peptide
36	BL	172	ARG	Sidechain
36	BL	174	ILE	Peptide
36	BL	33	LYS	Peptide
36	BL	38	ARG	Sidechain
36	BL	71	ARG	Sidechain
36	BL	81	TYR	Sidechain
37	BM	148	LYS	Mainchain,Peptide
37	BM	169	ALA	Mainchain
37	BM	84	LEU	Peptide
37	BM	85	ARG	Sidechain
38	BN	116	HIS	Peptide
38	BN	131	ARG	Sidechain
39	BO	72	LYS	Peptide
40	BP	110	ARG	Sidechain
40	BP	20	ARG	Sidechain
40	BP	38	ARG	Sidechain
40	BP	89	LEU	Peptide
42	BR	12	ARG	Sidechain
42	BR	32	ARG	Sidechain
42	BR	86	ARG	Sidechain
43	BS	47	ARG	Sidechain
44	BT	125	ARG	Sidechain
45	BU	63	LYS	Peptide
46	BV	112	LEU	Peptide
46	BV	12	ARG	Sidechain
46	BV	91	LYS	Peptide
47	BW	28	ARG	Sidechain
47	BW	74	ARG	Sidechain
48	BX	10	ARG	Sidechain

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Mol	Chain	Res	Type	Group
49	BY	47	TYR	Sidechain
50	BZ	32	LYS	Peptide
50	BZ	42	ARG	Sidechain
50	BZ	55	LYS	Peptide
50	BZ	56	PRO	Peptide
50	BZ	60	LYS	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	37458	0	18828	602	0
2	AB	1500	0	1524	1	0
3	AC	1469	0	1534	1	0
4	AD	1018	0	1067	1	0
5	AE	1207	0	1274	0	0
6	AG	1456	0	1523	2	0
7	AH	992	0	1027	39	0
8	AI	1087	0	1145	19	0
9	AJ	771	0	829	2	0
10	AK	924	0	949	0	0
11	AL	906	0	970	1	0
12	AM	1077	0	1107	19	0
13	AN	417	0	408	22	0
14	AO	694	0	740	13	0
15	AQ	643	0	688	2	0
16	AS	548	0	574	3	0
17	AR	2410	0	2367	538	0
18	AT	1102	0	1112	38	0
19	A7	1648	0	830	56	0
20	B0	881	0	926	1	0
21	B1	424	0	459	1	0
22	B2	752	0	803	1	0
23	B8	947	0	972	19	0
24	B9	539	0	555	1	0
25	BA	1683	0	1772	3	0
26	BB	1848	0	1908	16	0
27	BC	2887	0	2964	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
28	BD	1950	0	2019	3	0
29	BE	1913	0	1842	1	0
30	BF	1561	0	1633	1	0
31	BG	844	0	909	0	0
32	BH	1418	0	1483	4	0
33	BI	1326	0	1368	2	0
34	BJ	1195	0	1230	15	0
35	BK	1038	0	1110	23	0
36	BL	1618	0	1673	1	0
37	BM	1317	0	1407	63	0
38	BN	1189	0	1201	0	0
39	BO	931	0	1010	0	0
40	BP	1317	0	1389	113	0
41	BQ	893	0	924	0	0
42	BR	977	0	1026	15	0
43	BS	371	0	382	1	0
44	BT	642	0	679	0	0
45	BU	916	0	996	0	0
46	BV	1123	0	1160	2	0
47	BW	663	0	700	0	0
48	BX	605	0	661	1	0
49	BY	403	0	398	0	0
50	BZ	749	0	815	0	0
51	B3	2403	0	1219	0	0
52	B4	3329	0	1685	0	0
53	B5	67775	0	34047	227	0
All	All	165754	0	111821	1335	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 5.

All (1335) 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)
35:BK:121:PHE:CZ	35:BK:121:PHE:CE1	1.78	1.71
35:BK:121:PHE:CE2	35:BK:121:PHE:CZ	1.79	1.69
35:BK:121:PHE:CE2	35:BK:121:PHE:CD2	1.76	1.69
35:BK:121:PHE:CD1	35:BK:121:PHE:CE1	1.80	1.69
35:BK:121:PHE:CG	35:BK:121:PHE:CD2	1.82	1.68
35:BK:121:PHE:CD1	35:BK:121:PHE:CG	1.81	1.67
1:AA:1501:A:H1'	1:AA:1546:G:C2	1.33	1.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1756:U:H5'	53:B5:2262:A:C2	1.34	1.62
8:AI:96:TYR:CE1	8:AI:96:TYR:CD1	1.89	1.59
8:AI:96:TYR:CD2	8:AI:96:TYR:CE2	1.88	1.58
8:AI:96:TYR:CD2	8:AI:96:TYR:CG	1.90	1.58
8:AI:96:TYR:CE1	8:AI:96:TYR:CZ	1.91	1.57
8:AI:96:TYR:CG	8:AI:96:TYR:CD1	1.91	1.56
1:AA:1200:A:C5	1:AA:1200:A:C6	1.87	1.56
8:AI:96:TYR:CZ	8:AI:96:TYR:CE2	1.90	1.55
1:AA:1200:A:N3	1:AA:1200:A:C2	1.75	1.54
1:AA:1200:A:N1	1:AA:1200:A:C2	1.74	1.54
37:BM:111:PRO:HA	37:BM:163:ALA:CB	1.36	1.52
12:AM:110:ARG:NE	34:BJ:116:TYR:CD1	1.76	1.52
1:AA:702:G:C4	40:BP:172:ALA:O	1.64	1.50
1:AA:856:A:C6	1:AA:856:A:C5	1.99	1.50
37:BM:111:PRO:CA	37:BM:163:ALA:HB1	1.04	1.50
1:AA:702:G:C6	40:BP:172:ALA:HB1	0.98	1.49
1:AA:1200:A:C4	1:AA:1200:A:C5	1.86	1.48
1:AA:1630:C:H4'	1:AA:1636:G:N2	1.16	1.48
1:AA:1776:G:C4'	53:B5:2194:G:H5''	1.43	1.48
1:AA:1200:A:N1	1:AA:1200:A:C6	1.82	1.48
1:AA:702:G:C6	40:BP:172:ALA:CB	1.93	1.48
1:AA:699:U:H3'	40:BP:169:ALA:CB	1.42	1.47
1:AA:856:A:C2	1:AA:856:A:N1	1.82	1.47
1:AA:1200:A:N3	1:AA:1200:A:C4	1.82	1.47
37:BM:111:PRO:HB3	37:BM:163:ALA:CA	1.41	1.47
1:AA:1501:A:C6	1:AA:1546:G:C8	2.02	1.46
1:AA:1200:A:C4	13:AN:16:LYS:HA	1.46	1.45
1:AA:701:U:C4	40:BP:169:ALA:O	1.67	1.44
1:AA:856:A:C5	1:AA:856:A:C4	1.97	1.44
1:AA:501:U:C4	1:AA:501:U:C5	2.04	1.43
1:AA:856:A:C2	1:AA:856:A:N3	1.84	1.43
12:AM:110:ARG:NH2	34:BJ:116:TYR:CZ	1.82	1.42
1:AA:1501:A:N7	1:AA:1546:G:H2'	1.29	1.42
1:AA:702:G:C8	40:BP:175:ALA:HB3	1.54	1.40
1:AA:856:A:C6	1:AA:856:A:N1	1.87	1.40
1:AA:501:U:C6	1:AA:501:U:N1	1.89	1.40
1:AA:1756:U:H5'	53:B5:2262:A:N1	1.17	1.39
1:AA:501:U:C4	1:AA:501:U:N3	1.89	1.38
1:AA:856:A:N3	1:AA:856:A:C4	1.91	1.38
1:AA:501:U:N3	1:AA:501:U:C2	1.91	1.38
1:AA:1776:G:OP1	53:B5:2193:U:C5'	1.71	1.37

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1755:G:N2	53:B5:2262:A:O2'	1.57	1.36
1:AA:1757:C:O4'	53:B5:2263:C:C5'	1.74	1.35
1:AA:1501:A:C6	1:AA:1546:G:N7	1.93	1.35
1:AA:1200:A:C5	13:AN:16:LYS:HA	1.58	1.35
1:AA:501:U:C2	1:AA:501:U:N1	1.94	1.34
1:AA:1756:U:C5'	53:B5:2262:A:C2	2.09	1.33
37:BM:169:ALA:C	37:BM:171:ALA:N	1.78	1.33
1:AA:1501:A:H1'	1:AA:1546:G:N2	1.43	1.33
1:AA:702:G:O6	40:BP:172:ALA:CB	1.68	1.33
12:AM:110:ARG:NE	34:BJ:116:TYR:CE1	1.74	1.33
14:AO:149:LEU:HD12	53:B5:847:A:OP1	1.22	1.33
1:AA:1521:G:H5''	1:AA:1522:A:P	1.68	1.32
1:AA:1630:C:C4'	1:AA:1636:G:N2	1.90	1.32
1:AA:702:G:N3	40:BP:176:ALA:HB3	1.44	1.32
1:AA:1665:A:OP1	53:B5:1936:A:C5'	1.78	1.32
1:AA:1501:A:N1	1:AA:1546:G:N7	1.77	1.31
1:AA:699:U:C3'	40:BP:169:ALA:HB3	1.60	1.31
1:AA:1756:U:O3'	53:B5:2263:C:C4'	1.64	1.31
1:AA:1502:G:C5	1:AA:1547:C:C4	2.17	1.30
1:AA:1501:A:C1'	1:AA:1546:G:C2	2.11	1.30
1:AA:1655:U:C4	53:B5:2329:C:H1'	1.50	1.30
1:AA:972:A:H5'	53:B5:848:A:C2	1.64	1.29
37:BM:108:ILE:HG23	37:BM:160:ALA:CB	1.60	1.28
1:AA:1755:G:O2'	53:B5:2262:A:C2	1.84	1.27
1:AA:972:A:H5'	53:B5:848:A:N3	1.48	1.27
1:AA:1521:G:C2	18:AT:78:LYS:CG	2.16	1.27
1:AA:858:G:C2	7:AH:61:ILE:CB	2.16	1.27
1:AA:858:G:C6	7:AH:61:ILE:CB	2.18	1.27
37:BM:108:ILE:CG2	37:BM:160:ALA:CB	2.11	1.27
1:AA:1735:G:H4'	53:B5:1933:A:O2'	1.31	1.27
1:AA:858:G:N1	7:AH:61:ILE:CB	1.99	1.26
1:AA:1501:A:C8	1:AA:1546:G:H2'	1.70	1.26
1:AA:1630:C:C4'	1:AA:1636:G:H21	1.46	1.26
1:AA:1501:A:C4	1:AA:1546:G:C4	2.16	1.26
1:AA:1776:G:OP1	53:B5:2193:U:H5''	1.14	1.25
1:AA:1521:G:C2	18:AT:78:LYS:HG3	1.68	1.25
1:AA:698:U:N1	40:BP:164:ALA:O	1.58	1.25
37:BM:111:PRO:CA	37:BM:163:ALA:CB	2.01	1.24
1:AA:698:U:C2	40:BP:164:ALA:O	1.90	1.24
1:AA:699:U:C3'	40:BP:169:ALA:CB	2.13	1.24
12:AM:110:ARG:NH2	34:BJ:116:TYR:CE2	2.04	1.23

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
14:AO:123:HIS:ND1	53:B5:846:A:N3	1.84	1.23
1:AA:1502:G:N7	1:AA:1547:C:C2	2.07	1.23
1:AA:1776:G:H5''	53:B5:2194:G:C5'	1.69	1.22
1:AA:702:G:C5	40:BP:172:ALA:HB1	1.75	1.22
1:AA:698:U:C6	40:BP:164:ALA:O	1.91	1.22
1:AA:1574:A:N3	19:A7:41:U:O4'	1.70	1.21
1:AA:971:G:H21	53:B5:846:A:N6	1.37	1.21
1:AA:1574:A:N3	19:A7:41:U:C4'	1.84	1.21
1:AA:1653:A:C6	53:B5:2302:G:C5'	2.16	1.20
1:AA:1657:A:H5''	42:BR:67:PRO:CD	1.72	1.20
1:AA:702:G:O6	40:BP:172:ALA:HB1	1.02	1.19
37:BM:111:PRO:CB	37:BM:163:ALA:HB1	1.72	1.19
1:AA:1776:G:O3'	53:B5:2194:G:C4'	1.90	1.19
1:AA:1756:U:C1'	53:B5:2262:A:H2'	1.72	1.19
1:AA:799:A:N3	40:BP:161:ALA:CB	1.99	1.19
1:AA:1756:U:C3'	53:B5:2263:C:H4'	1.52	1.18
1:AA:1635:C:C4'	1:AA:1636:G:OP2	1.84	1.18
1:AA:1757:C:O4'	53:B5:2263:C:H5'	1.31	1.17
1:AA:1461:C:OP1	19:A7:30:G:H1'	1.40	1.17
1:AA:1776:G:O3'	53:B5:2194:G:H4'	0.99	1.17
1:AA:1630:C:C5'	1:AA:1636:G:H21	1.56	1.16
1:AA:1756:U:H1'	53:B5:2262:A:C2'	1.75	1.15
1:AA:921:G:H4'	26:BB:141:PRO:HD3	1.28	1.15
1:AA:1776:G:H4'	53:B5:2194:G:C5'	1.76	1.15
17:AR:34:LEU:HD21	17:AR:42:LEU:HD21	1.27	1.15
1:AA:1635:C:H4'	1:AA:1636:G:OP2	1.43	1.15
1:AA:1521:G:N1	18:AT:78:LYS:CG	2.01	1.15
17:AR:34:LEU:HG	17:AR:42:LEU:HD11	1.24	1.15
1:AA:1655:U:C4	53:B5:2329:C:C1'	2.26	1.15
1:AA:1776:G:C5'	53:B5:2194:G:C5'	2.23	1.14
1:AA:1121:A:C1'	53:B5:2191:U:C5'	2.24	1.14
17:AR:38:ARG:HA	17:AR:67:ILE:HG23	1.29	1.14
37:BM:111:PRO:CB	37:BM:163:ALA:CA	2.26	1.14
1:AA:1521:G:H1'	18:AT:86:ARG:NH2	1.63	1.14
1:AA:1121:A:H1'	53:B5:2191:U:H5''	1.16	1.14
1:AA:1502:G:C5	1:AA:1547:C:N3	2.14	1.13
1:AA:1653:A:H1'	53:B5:2302:G:OP2	1.45	1.13
1:AA:1521:G:N2	18:AT:78:LYS:CD	2.10	1.13
1:AA:1521:G:O2'	18:AT:83:ALA:HA	1.48	1.13
1:AA:1502:G:C8	1:AA:1547:C:N3	2.17	1.13
1:AA:1502:G:C6	1:AA:1547:C:C4	2.35	1.13

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
37:BM:108:ILE:CG2	37:BM:160:ALA:HB1	1.72	1.13
1:AA:1521:G:N1	18:AT:78:LYS:HG3	1.28	1.13
1:AA:1200:A:C5	13:AN:16:LYS:CA	2.32	1.12
1:AA:1521:G:C5	18:AT:79:LEU:N	2.17	1.12
1:AA:1200:A:C4	13:AN:16:LYS:CA	2.32	1.12
1:AA:1502:G:N7	1:AA:1547:C:N3	1.97	1.12
1:AA:1501:A:C5	1:AA:1546:G:C4	2.36	1.12
1:AA:1757:C:O4'	53:B5:2263:C:H5''	1.50	1.12
1:AA:1200:A:C2	13:AN:16:LYS:CA	2.33	1.12
1:AA:1630:C:H4'	1:AA:1636:G:C2	1.83	1.12
1:AA:1776:G:C5'	53:B5:2194:G:H5'	1.78	1.12
37:BM:108:ILE:CG2	37:BM:160:ALA:HB2	1.76	1.12
17:AR:209:THR:HG23	17:AR:210:LEU:HD22	1.16	1.12
1:AA:1756:U:C5'	53:B5:2262:A:N1	2.07	1.12
1:AA:1776:G:C4'	53:B5:2194:G:C5'	2.26	1.12
1:AA:1665:A:OP1	53:B5:1936:A:H5'	0.94	1.12
1:AA:1778:G:OP1	53:B5:2274:U:H5'	1.44	1.11
1:AA:858:G:C5	7:AH:61:ILE:CB	2.31	1.11
37:BM:151:ALA:O	37:BM:155:ALA:CB	1.98	1.11
14:AO:149:LEU:CD1	53:B5:847:A:OP1	1.98	1.10
1:AA:1653:A:C1'	53:B5:2302:G:OP2	1.96	1.10
14:AO:149:LEU:HD13	53:B5:847:A:H5'	1.21	1.10
1:AA:1521:G:H5''	1:AA:1522:A:OP2	1.47	1.10
1:AA:1665:A:O4'	53:B5:1935:G:H5'	1.52	1.09
1:AA:1657:A:C5'	42:BR:67:PRO:HG2	1.81	1.09
1:AA:1574:A:H2'	19:A7:40:5MC:O2'	1.41	1.09
1:AA:699:U:O2	40:BP:170:ALA:HB3	1.35	1.09
37:BM:169:ALA:O	37:BM:170:ALA:C	1.79	1.09
1:AA:1501:A:N7	1:AA:1546:G:C2'	2.15	1.08
1:AA:1756:U:C3'	53:B5:2263:C:C4'	2.16	1.08
1:AA:1664:U:O3'	53:B5:1935:G:O2'	1.71	1.08
17:AR:211:ILE:HD11	17:AR:225:LEU:HD13	1.16	1.08
1:AA:1574:A:C2	19:A7:41:U:O4'	2.05	1.08
1:AA:1200:A:C6	13:AN:16:LYS:CA	2.37	1.08
1:AA:856:A:C5	7:AH:33:VAL:N	2.22	1.08
1:AA:1632:C:H4'	1:AA:1633:A:OP2	1.50	1.08
1:AA:1776:G:H5'	53:B5:2194:G:OP1	1.54	1.07
1:AA:1521:G:C5'	1:AA:1522:A:OP2	2.03	1.07
1:AA:1521:G:C2	18:AT:78:LYS:CB	2.38	1.07
1:AA:972:A:C5'	53:B5:848:A:N3	2.16	1.07
1:AA:1521:G:C5'	1:AA:1522:A:P	2.33	1.06

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
17:AR:164:ASP:HB2	17:AR:168:THR:HA	1.09	1.06
1:AA:856:A:C4	7:AH:32:LYS:C	2.29	1.06
1:AA:1501:A:C5	1:AA:1546:G:C8	2.44	1.06
37:BM:108:ILE:HG23	37:BM:160:ALA:HB1	1.32	1.06
14:AO:149:LEU:HD12	53:B5:847:A:P	1.95	1.06
37:BM:111:PRO:CB	37:BM:163:ALA:CB	2.28	1.06
1:AA:1757:C:C4'	53:B5:2263:C:H5''	1.78	1.05
1:AA:702:G:N7	40:BP:172:ALA:HA	1.71	1.05
1:AA:699:U:O2	40:BP:170:ALA:CB	1.77	1.05
1:AA:699:U:C1'	40:BP:169:ALA:HB3	1.77	1.05
1:AA:800:U:OP1	40:BP:160:ALA:HB1	1.56	1.05
1:AA:858:G:C4	7:AH:61:ILE:CB	2.39	1.05
17:AR:260:ILE:HG13	17:AR:284:ALA:HB1	1.38	1.05
1:AA:1657:A:H5'	42:BR:67:PRO:HG2	1.33	1.05
17:AR:258:THR:HG21	17:AR:261:LYS:HE2	1.39	1.05
17:AR:164:ASP:HB2	17:AR:168:THR:CA	1.87	1.05
1:AA:856:A:C4	7:AH:33:VAL:N	2.24	1.05
1:AA:856:A:C5	7:AH:32:LYS:C	2.30	1.05
1:AA:1744:A:C2	53:B5:2302:G:O5'	2.09	1.05
23:B8:43:LYS:NZ	35:BK:121:PHE:CE1	2.25	1.04
1:AA:1200:A:C5	13:AN:16:LYS:N	2.25	1.04
12:AM:110:ARG:CZ	34:BJ:116:TYR:CE1	2.40	1.04
1:AA:1734:G:H21	53:B5:1934:G:H4'	1.19	1.04
1:AA:856:A:C6	7:AH:33:VAL:N	2.26	1.04
37:BM:111:PRO:CB	37:BM:163:ALA:HA	1.86	1.04
14:AO:149:LEU:HD13	53:B5:847:A:C5'	1.87	1.04
1:AA:702:G:C5	40:BP:172:ALA:O	2.10	1.03
1:AA:1200:A:C4	13:AN:16:LYS:N	2.25	1.03
1:AA:799:A:C4'	40:BP:161:ALA:O	2.04	1.03
1:AA:856:A:C2	7:AH:33:VAL:N	2.26	1.03
8:AI:93:HIS:CE1	17:AR:59:ARG:HH22	1.76	1.03
17:AR:257:ALA:HA	17:AR:283:LYS:HD2	1.07	1.03
17:AR:83:ALA:HB2	17:AR:113:VAL:HG13	1.41	1.03
1:AA:1501:A:N6	1:AA:1546:G:C8	2.26	1.03
1:AA:699:U:O4'	40:BP:166:ALA:HA	1.56	1.03
1:AA:1744:A:H4'	53:B5:2291:A:H1'	1.37	1.03
1:AA:922:A:H5'	26:BB:109:GLU:OE1	1.57	1.03
37:BM:111:PRO:HB3	37:BM:163:ALA:HA	1.07	1.03
1:AA:856:A:C2	7:AH:32:LYS:C	2.32	1.03
1:AA:1630:C:C5'	1:AA:1636:G:N2	2.18	1.03
17:AR:93:ASP:HB3	17:AR:96:THR:HG22	1.41	1.03

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1200:A:C2	13:AN:16:LYS:N	2.27	1.02
1:AA:699:U:C2'	40:BP:169:ALA:HB3	1.89	1.02
1:AA:699:U:H3'	40:BP:169:ALA:HB1	1.04	1.02
1:AA:701:U:C5	40:BP:169:ALA:O	2.11	1.02
1:AA:702:G:C8	40:BP:175:ALA:CB	2.41	1.02
1:AA:1521:G:N2	18:AT:78:LYS:HD2	1.73	1.02
19:A7:71:G:H5'	53:B5:2236:G:H4'	1.42	1.02
1:AA:702:G:C5	40:BP:172:ALA:CA	2.43	1.02
17:AR:286:GLU:HG2	17:AR:289:ALA:HA	1.41	1.02
1:AA:1121:A:H1'	53:B5:2191:U:C5'	1.85	1.02
1:AA:799:A:N3	40:BP:161:ALA:HB2	1.43	1.01
1:AA:1520:U:H5''	1:AA:1522:A:OP2	1.59	1.01
1:AA:1653:A:C6	53:B5:2302:G:H5''	1.88	1.01
17:AR:210:LEU:HD12	17:AR:222:LEU:HD21	1.40	1.01
1:AA:1000:U:OP1	19:A7:38:A:H5'	1.58	1.01
1:AA:799:A:O4'	40:BP:161:ALA:O	1.71	1.01
1:AA:699:U:O4'	40:BP:166:ALA:CA	2.07	1.01
1:AA:1574:A:C2	19:A7:41:U:C4'	2.42	1.01
1:AA:1755:G:O2'	53:B5:2262:A:H2	1.25	1.00
1:AA:1756:U:H1'	53:B5:2262:A:H2'	1.01	1.00
1:AA:1200:A:C6	13:AN:16:LYS:N	2.29	1.00
1:AA:1657:A:C5'	42:BR:67:PRO:CG	2.39	1.00
1:AA:1501:A:C5	1:AA:1546:G:C5	2.44	1.00
1:AA:1756:U:O4'	53:B5:2262:A:N3	1.93	1.00
1:AA:856:A:C6	7:AH:32:LYS:C	2.35	1.00
1:AA:1777:U:P	53:B5:2194:G:H4'	2.02	1.00
1:AA:1744:A:C4'	53:B5:2291:A:H1'	1.92	0.99
1:AA:1121:A:C4'	53:B5:2191:U:H5'	1.92	0.99
1:AA:912:G:O6	53:B5:2207:A:O5'	1.71	0.99
17:AR:12:THR:HG22	17:AR:311:ARG:HG2	1.40	0.99
1:AA:1635:C:C5'	1:AA:1636:G:OP2	2.10	0.99
1:AA:1574:A:C2'	19:A7:40:5MC:O2'	2.11	0.99
1:AA:1501:A:C1'	1:AA:1546:G:N2	2.15	0.99
1:AA:702:G:H1'	40:BP:175:ALA:CB	1.78	0.99
17:AR:259:GLY:HA2	17:AR:284:ALA:HA	1.41	0.99
1:AA:1756:U:C4'	53:B5:2262:A:C2	2.46	0.98
1:AA:1502:G:C6	1:AA:1547:C:C5	2.52	0.98
1:AA:1502:G:C8	1:AA:1547:C:C2	2.50	0.98
1:AA:1657:A:H5'	42:BR:67:PRO:CG	1.93	0.98
1:AA:1121:A:C1'	53:B5:2191:U:H5'	1.91	0.98
1:AA:993:G:H5''	53:B5:2195:C:OP2	1.64	0.98

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1756:U:O3'	53:B5:2263:C:C1'	1.98	0.97
1:AA:1521:G:N2	18:AT:78:LYS:NZ	2.11	0.97
1:AA:1460:G:H4'	19:A7:30:G:H5'	1.46	0.97
1:AA:1756:U:C2'	53:B5:2263:C:OP1	2.12	0.97
1:AA:1653:A:H2	53:B5:2301:U:HO2'	1.02	0.97
1:AA:1756:U:O2'	53:B5:2263:C:O5'	1.80	0.97
1:AA:971:G:N2	53:B5:846:A:N6	2.12	0.96
12:AM:110:ARG:CZ	34:BJ:116:TYR:CZ	2.48	0.96
1:AA:972:A:C5'	53:B5:848:A:C2	2.49	0.96
1:AA:1655:U:O2	53:B5:2124:G:N2	1.99	0.96
1:AA:701:U:O4	40:BP:169:ALA:O	1.80	0.96
1:AA:699:U:C2	40:BP:167:ALA:O	2.20	0.95
17:AR:40:LYS:HE2	17:AR:66:HIS:HA	1.48	0.95
1:AA:699:U:C2	40:BP:167:ALA:C	2.39	0.95
37:BM:111:PRO:HB3	37:BM:163:ALA:C	1.86	0.95
37:BM:108:ILE:HG22	37:BM:160:ALA:HB2	1.46	0.95
1:AA:1502:G:C4	1:AA:1547:C:N4	2.35	0.94
17:AR:286:GLU:HG3	17:AR:305:TYR:CG	2.02	0.94
17:AR:34:LEU:CG	17:AR:42:LEU:HD11	1.97	0.94
37:BM:151:ALA:O	37:BM:155:ALA:HB3	1.66	0.94
17:AR:164:ASP:CB	17:AR:168:THR:HA	1.97	0.94
37:BM:169:ALA:C	37:BM:171:ALA:H	1.68	0.94
17:AR:257:ALA:CA	17:AR:283:LYS:HD2	1.97	0.93
1:AA:1776:G:C5'	53:B5:2194:G:H5''	1.92	0.93
1:AA:1735:G:C4'	53:B5:1933:A:O2'	2.16	0.93
19:A7:63:C:H5''	25:BA:45:ARG:HD2	1.49	0.93
1:AA:1520:U:H4'	1:AA:1522:A:C8	2.03	0.93
1:AA:1657:A:H5''	42:BR:67:PRO:HD2	1.48	0.93
1:AA:865:A:H8	1:AA:865:A:OP1	1.51	0.93
1:AA:1776:G:H5''	53:B5:2194:G:H5'	0.94	0.93
1:AA:1735:G:H4'	53:B5:1933:A:HO2'	1.31	0.93
1:AA:1521:G:C2	18:AT:78:LYS:HB2	2.02	0.93
17:AR:46:LYS:HB2	17:AR:58:VAL:CG1	1.98	0.93
37:BM:155:ALA:O	37:BM:158:ALA:N	2.01	0.92
17:AR:164:ASP:HB3	17:AR:183:LEU:H	1.33	0.92
17:AR:36:ALA:HB1	17:AR:68:VAL:HG13	1.46	0.92
1:AA:993:G:C5'	53:B5:2195:C:OP2	2.17	0.92
1:AA:1501:A:C5	1:AA:1546:G:N9	2.37	0.92
1:AA:702:G:N9	40:BP:175:ALA:HB3	1.84	0.92
12:AM:110:ARG:CD	34:BJ:116:TYR:CD1	2.53	0.92
1:AA:702:G:N3	40:BP:176:ALA:CB	2.30	0.92

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
17:AR:156:VAL:HB	17:AR:167:VAL:CG2	2.00	0.92
1:AA:1655:U:C5	53:B5:2329:C:H1'	2.05	0.92
17:AR:83:ALA:HB2	17:AR:113:VAL:CG1	1.99	0.92
23:B8:43:LYS:NZ	35:BK:121:PHE:CD1	2.38	0.92
1:AA:1501:A:C6	1:AA:1546:G:C5	2.57	0.92
1:AA:1756:U:C1'	53:B5:2263:C:OP1	2.18	0.92
1:AA:1755:G:C2'	53:B5:2262:A:H2	1.83	0.91
1:AA:1501:A:C2	1:AA:1546:G:N7	2.25	0.91
1:AA:1657:A:C5'	42:BR:67:PRO:CD	2.49	0.91
1:AA:1744:A:C5	53:B5:2303:A:OP2	2.24	0.91
1:AA:1460:G:O3'	19:A7:30:G:O4'	1.87	0.91
17:AR:258:THR:CG2	17:AR:261:LYS:HE2	2.00	0.91
17:AR:286:GLU:HB3	17:AR:288:HIS:H	1.36	0.91
19:A7:56:C:C5	53:B5:2484:A:H1'	2.06	0.91
1:AA:1521:G:H22	18:AT:78:LYS:NZ	1.69	0.90
1:AA:1756:U:O3'	53:B5:2263:C:H4'	1.42	0.90
17:AR:259:GLY:HA2	17:AR:284:ALA:CA	2.00	0.90
14:AO:149:LEU:CD1	53:B5:847:A:P	2.58	0.90
8:AI:93:HIS:CE1	17:AR:59:ARG:NH2	2.40	0.90
1:AA:1460:G:C4'	19:A7:30:G:H5'	2.02	0.90
23:B8:43:LYS:CE	35:BK:121:PHE:CD1	2.55	0.90
1:AA:1744:A:N7	53:B5:2303:A:OP2	1.89	0.90
1:AA:1756:U:O3'	53:B5:2263:C:O4'	1.89	0.89
17:AR:283:LYS:HD3	17:AR:288:HIS:HA	1.54	0.89
17:AR:61:PHE:CE2	17:AR:97:GLY:HA2	2.06	0.89
8:AI:93:HIS:HE1	17:AR:59:ARG:HH22	1.09	0.89
17:AR:266:ASP:HB3	17:AR:267:PRO:HD3	1.52	0.89
1:AA:1521:G:H1	18:AT:78:LYS:HG3	1.07	0.89
1:AA:800:U:OP1	40:BP:160:ALA:CB	2.21	0.89
1:AA:701:U:O2	40:BP:174:ALA:N	2.04	0.88
17:AR:93:ASP:HB3	17:AR:96:THR:CG2	2.03	0.88
1:AA:972:A:O4'	53:B5:847:A:H2	1.56	0.88
1:AA:1757:C:OP1	53:B5:2264:U:C5'	2.21	0.88
1:AA:972:A:O4'	53:B5:847:A:C2	2.25	0.88
1:AA:1520:U:H4'	1:AA:1522:A:N7	1.86	0.88
17:AR:257:ALA:HA	17:AR:283:LYS:CD	2.01	0.88
1:AA:1632:C:C4'	1:AA:1633:A:OP2	2.21	0.88
1:AA:1200:A:N1	13:AN:16:LYS:CA	2.37	0.88
1:AA:1502:G:C4	1:AA:1547:C:C4	2.62	0.88
1:AA:922:A:C5'	26:BB:109:GLU:OE1	2.21	0.88
1:AA:1121:A:O4'	53:B5:2191:U:H5'	1.73	0.88

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1756:U:H2'	53:B5:2263:C:C5'	2.02	0.88
37:BM:151:ALA:O	37:BM:155:ALA:HB2	1.73	0.88
1:AA:1521:G:H1'	18:AT:86:ARG:HH22	1.35	0.88
1:AA:702:G:C5	40:BP:172:ALA:C	2.47	0.88
1:AA:1777:U:OP1	53:B5:2194:G:H1'	1.72	0.88
1:AA:1521:G:H22	18:AT:78:LYS:HZ2	1.18	0.88
17:AR:283:LYS:HD3	17:AR:288:HIS:CA	2.04	0.88
17:AR:286:GLU:HG2	17:AR:289:ALA:CA	2.04	0.88
17:AR:38:ARG:HA	17:AR:67:ILE:CG2	2.03	0.88
23:B8:43:LYS:HB2	35:BK:121:PHE:CD1	2.10	0.87
1:AA:1121:A:C1'	53:B5:2191:U:H5''	1.77	0.87
1:AA:1501:A:N9	1:AA:1546:G:N3	2.23	0.87
17:AR:260:ILE:HG13	17:AR:284:ALA:CB	2.03	0.87
17:AR:197:SER:CB	17:AR:216:LYS:HB3	2.04	0.87
1:AA:1521:G:N2	18:AT:78:LYS:CG	2.34	0.87
17:AR:89:LEU:HD21	17:AR:124:SER:HB3	1.53	0.87
1:AA:856:A:C5	7:AH:32:LYS:HA	2.08	0.87
1:AA:1521:G:N2	18:AT:78:LYS:HG3	1.90	0.87
37:BM:108:ILE:HG23	37:BM:160:ALA:HB2	1.42	0.87
1:AA:1756:U:H2'	53:B5:2263:C:H5'	1.57	0.86
23:B8:43:LYS:HE3	35:BK:121:PHE:HD1	1.40	0.86
23:B8:43:LYS:CG	35:BK:121:PHE:CE1	2.58	0.86
17:AR:283:LYS:CB	17:AR:288:HIS:HA	2.05	0.86
1:AA:1757:C:OP1	53:B5:2264:U:H5'	1.76	0.86
17:AR:52:GLN:HA	17:AR:52:GLN:HE21	1.41	0.86
1:AA:702:G:C5	40:BP:172:ALA:CB	2.41	0.86
1:AA:922:A:OP1	26:BB:109:GLU:OE2	1.93	0.86
1:AA:865:A:C8	1:AA:865:A:OP1	2.28	0.86
1:AA:1744:A:N3	53:B5:2302:G:O5'	2.04	0.86
1:AA:1516:C:H6	1:AA:1516:C:OP2	1.57	0.86
37:BM:111:PRO:HB3	37:BM:163:ALA:CB	1.99	0.86
1:AA:1776:G:OP1	53:B5:2193:U:H5'	1.72	0.85
17:AR:164:ASP:O	17:AR:183:LEU:HB2	1.75	0.85
1:AA:701:U:C2'	40:BP:176:ALA:O	2.23	0.85
1:AA:699:U:H3	40:BP:171:ALA:H	1.24	0.85
1:AA:1574:A:C2	19:A7:41:U:H4'	2.08	0.85
12:AM:110:ARG:CZ	34:BJ:116:TYR:CD1	2.57	0.85
17:AR:281:TYR:O	17:AR:287:PRO:HA	1.76	0.85
17:AR:283:LYS:CG	17:AR:288:HIS:HA	2.06	0.85
17:AR:165:ASP:HB2	17:AR:183:LEU:O	1.77	0.85
1:AA:1501:A:N9	1:AA:1546:G:C2	2.45	0.85

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1653:A:C5	53:B5:2302:G:H5''	2.08	0.85
1:AA:1000:U:OP1	19:A7:38:A:C5'	2.24	0.85
17:AR:164:ASP:HB3	17:AR:183:LEU:N	1.91	0.84
1:AA:1777:U:OP1	53:B5:2194:G:C1'	2.24	0.84
17:AR:209:THR:HG23	17:AR:210:LEU:CD2	2.05	0.84
17:AR:29:GLN:HB2	17:AR:296:ALA:HB3	1.57	0.84
1:AA:1734:G:N2	53:B5:1934:G:H4'	1.91	0.84
17:AR:193:ILE:H	17:AR:193:ILE:HD13	1.42	0.84
17:AR:164:ASP:CB	17:AR:183:LEU:H	1.90	0.84
37:BM:150:ALA:O	37:BM:154:ALA:N	2.10	0.83
1:AA:702:G:H8	40:BP:175:ALA:HB3	1.36	0.83
17:AR:288:HIS:HB3	17:AR:306:THR:CG2	2.09	0.83
17:AR:33:LEU:CD2	17:AR:45:TRP:HB2	2.08	0.83
17:AR:164:ASP:CG	17:AR:182:ASN:HA	1.98	0.83
17:AR:197:SER:HB3	17:AR:216:LYS:HB3	1.59	0.83
1:AA:858:G:N3	7:AH:61:ILE:CB	2.41	0.83
17:AR:169:ILE:HG13	17:AR:183:LEU:HD21	1.60	0.83
1:AA:699:U:C4'	40:BP:169:ALA:HB3	2.09	0.83
19:A7:71:G:H5'	53:B5:2236:G:C4'	2.08	0.83
1:AA:1200:A:N3	13:AN:16:LYS:CA	2.42	0.82
1:AA:1653:A:H2	53:B5:2301:U:O2'	1.61	0.82
17:AR:262:VAL:HG23	17:AR:271:VAL:HG12	1.61	0.82
17:AR:133:VAL:HG13	17:AR:141:LEU:HB2	1.60	0.82
1:AA:1339:C:H5'	17:AR:102:ARG:NH2	1.94	0.82
1:AA:699:U:N3	40:BP:167:ALA:O	2.13	0.82
1:AA:921:G:H1'	26:BB:141:PRO:HG3	1.59	0.82
23:B8:43:LYS:HE3	35:BK:121:PHE:CD1	2.14	0.82
17:AR:21:THR:HG21	17:AR:68:VAL:O	1.79	0.82
17:AR:163:ASP:HB2	17:AR:165:ASP:O	1.80	0.82
17:AR:203:THR:HG22	17:AR:212:ALA:HB3	1.62	0.82
1:AA:799:A:C4'	40:BP:161:ALA:C	2.47	0.82
17:AR:276:PRO:N	17:AR:285:ALA:HA	1.95	0.82
12:AM:110:ARG:HE	34:BJ:116:TYR:HE1	1.18	0.81
17:AR:305:TYR:CZ	17:AR:311:ARG:HB2	2.15	0.81
1:AA:1665:A:O4'	53:B5:1935:G:C5'	2.28	0.81
17:AR:177:MET:SD	17:AR:191:ASP:HB2	2.20	0.81
17:AR:34:LEU:CD2	17:AR:42:LEU:HD21	2.10	0.81
1:AA:1573:G:N1	19:A7:41:U:O2'	1.93	0.81
1:AA:699:U:H2'	40:BP:169:ALA:C	2.00	0.81
1:AA:702:G:C1'	40:BP:175:ALA:CB	2.58	0.81
17:AR:38:ARG:HG2	17:AR:67:ILE:HG21	1.62	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:699:U:H2'	40:BP:170:ALA:N	1.95	0.80
37:BM:148:LYS:HB2	37:BM:149:ALA:CB	2.11	0.80
1:AA:1501:A:C8	1:AA:1546:G:C2'	2.62	0.80
1:AA:702:G:N9	40:BP:172:ALA:O	2.14	0.80
17:AR:164:ASP:HB3	17:AR:182:ASN:HA	1.62	0.80
1:AA:702:G:N7	40:BP:172:ALA:CA	2.42	0.80
1:AA:1501:A:H1'	1:AA:1546:G:N1	1.95	0.80
1:AA:1521:G:N3	18:AT:78:LYS:HB2	1.90	0.80
17:AR:283:LYS:CD	17:AR:288:HIS:HA	2.10	0.80
17:AR:164:ASP:CG	17:AR:168:THR:HG22	2.02	0.80
19:A7:56:C:C6	53:B5:2484:A:H1'	2.17	0.80
1:AA:1776:G:C5'	53:B5:2194:G:OP1	2.30	0.80
1:AA:1665:A:P	53:B5:1935:G:O2'	2.39	0.80
1:AA:700:C:C6	40:BP:173:ALA:HB3	2.17	0.80
17:AR:283:LYS:HD3	17:AR:288:HIS:CG	2.16	0.80
1:AA:1655:U:O4	53:B5:2328:U:O2	1.99	0.80
17:AR:96:THR:HG23	17:AR:98:GLU:H	1.45	0.80
1:AA:856:A:C5	7:AH:32:LYS:CA	2.66	0.79
1:AA:1653:A:C2'	53:B5:2302:G:OP2	2.29	0.79
17:AR:164:ASP:CB	17:AR:182:ASN:HA	2.11	0.79
1:AA:699:U:C2'	40:BP:169:ALA:CB	2.56	0.79
1:AA:856:A:N1	7:AH:33:VAL:N	2.31	0.79
1:AA:856:A:N1	7:AH:32:LYS:C	2.36	0.79
1:AA:1521:G:O2'	18:AT:83:ALA:CA	2.30	0.79
17:AR:285:ALA:HB3	17:AR:313:TRP:CZ2	2.17	0.79
1:AA:921:G:C4'	26:BB:141:PRO:HD3	2.09	0.79
23:B8:43:LYS:CD	35:BK:121:PHE:CE1	2.66	0.79
1:AA:1657:A:H5''	42:BR:67:PRO:HD3	1.64	0.79
17:AR:115:ILE:CG2	17:AR:119:ALA:HA	2.13	0.78
1:AA:1775:G:O3'	53:B5:2193:U:OP1	2.00	0.78
1:AA:1776:G:P	53:B5:2193:U:H5''	2.23	0.78
1:AA:1573:G:H1	19:A7:41:U:HO2'	0.80	0.78
17:AR:83:ALA:HB1	17:AR:110:VAL:CG1	2.12	0.78
37:BM:168:ALA:C	37:BM:171:ALA:HB3	2.04	0.78
1:AA:1521:G:N2	18:AT:78:LYS:CE	2.46	0.78
1:AA:1516:C:C6	1:AA:1516:C:OP2	2.36	0.78
17:AR:211:ILE:HD11	17:AR:225:LEU:CD1	2.09	0.78
1:AA:699:U:C3'	40:BP:169:ALA:HB1	1.94	0.77
17:AR:13:LEU:HD11	17:AR:54:PHE:HB3	1.64	0.77
1:AA:1744:A:H4'	53:B5:2291:A:C1'	2.13	0.77
1:AA:1756:U:O2'	53:B5:2263:C:P	2.42	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1200:A:N3	13:AN:16:LYS:N	2.33	0.77
8:AI:97:VAL:HG21	17:AR:59:ARG:CZ	2.14	0.77
37:BM:108:ILE:HG22	37:BM:160:ALA:CB	2.03	0.77
1:AA:1745:G:O3'	53:B5:2303:A:H3'	1.81	0.77
19:A7:63:C:H5''	25:BA:45:ARG:CD	2.14	0.77
1:AA:856:A:N3	7:AH:33:VAL:N	2.33	0.77
1:AA:1521:G:H5''	1:AA:1522:A:OP1	1.82	0.77
1:AA:1736:U:H5'	53:B5:1932:A:H61	1.48	0.77
1:AA:700:C:C5	40:BP:169:ALA:O	2.36	0.77
37:BM:111:PRO:HA	37:BM:163:ALA:HB3	1.60	0.76
1:AA:1521:G:N2	18:AT:78:LYS:HZ2	1.78	0.76
1:AA:1502:G:C2	1:AA:1547:C:N4	2.54	0.76
1:AA:1573:G:C4	19:A7:42:G:OP1	2.34	0.76
17:AR:286:GLU:HA	17:AR:305:TYR:CD2	2.19	0.76
17:AR:286:GLU:HG3	17:AR:305:TYR:CD2	2.20	0.76
1:AA:1501:A:N9	1:AA:1546:G:C4	2.53	0.76
1:AA:702:G:C1'	40:BP:175:ALA:HB3	2.16	0.76
1:AA:856:A:N3	7:AH:32:LYS:C	2.39	0.76
17:AR:286:GLU:HB3	17:AR:288:HIS:N	2.00	0.76
17:AR:173:GLY:H	17:AR:199:ILE:CG2	1.99	0.76
1:AA:702:G:N3	40:BP:172:ALA:O	2.18	0.76
17:AR:170:ILE:HG22	17:AR:202:LEU:HD13	1.68	0.76
1:AA:1200:A:N1	13:AN:16:LYS:N	2.33	0.76
1:AA:1338:A:O3'	17:AR:102:ARG:CZ	2.33	0.76
17:AR:164:ASP:CA	17:AR:183:LEU:H	1.98	0.75
8:AI:97:VAL:HG21	17:AR:59:ARG:NH2	1.99	0.75
17:AR:19:TRP:CB	17:AR:38:ARG:HD2	2.16	0.75
1:AA:1502:G:C4	1:AA:1547:C:N3	2.54	0.75
1:AA:1521:G:H1'	18:AT:86:ARG:HH21	1.52	0.75
17:AR:172:ALA:CB	17:AR:199:ILE:HD12	2.16	0.75
1:AA:1000:U:H5'	19:A7:38:A:OP1	1.86	0.75
1:AA:1520:U:C4'	1:AA:1522:A:N7	2.49	0.75
17:AR:305:TYR:OH	17:AR:311:ARG:HB2	1.86	0.75
17:AR:288:HIS:HB3	17:AR:306:THR:HG23	1.69	0.75
17:AR:164:ASP:HB3	17:AR:182:ASN:CA	2.16	0.75
17:AR:33:LEU:HD23	17:AR:45:TRP:HB2	1.66	0.75
17:AR:13:LEU:CD1	17:AR:54:PHE:HB3	2.16	0.75
1:AA:701:U:O4	40:BP:172:ALA:CB	2.26	0.75
1:AA:1776:G:H4'	53:B5:2194:G:H5''	0.79	0.75
17:AR:86:ASP:OD1	17:AR:88:THR:HG22	1.87	0.75
1:AA:1756:U:C4'	53:B5:2262:A:N3	2.49	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1635:C:H5'	1:AA:1636:G:OP2	1.87	0.75
1:AA:1519:G:H4'	1:AA:1520:U:OP1	1.87	0.75
1:AA:1657:A:C4'	42:BR:67:PRO:HG2	2.17	0.75
1:AA:701:U:O4	40:BP:169:ALA:HA	1.87	0.74
1:AA:1502:G:C5	1:AA:1547:C:N4	2.53	0.74
1:AA:1521:G:C2	18:AT:78:LYS:CD	2.66	0.74
17:AR:114:ASP:OD1	17:AR:123:ILE:HG13	1.87	0.74
17:AR:126:SER:HB3	17:AR:128:ASP:OD1	1.87	0.74
17:AR:173:GLY:O	17:AR:199:ILE:HG22	1.87	0.74
8:AI:93:HIS:HE1	17:AR:59:ARG:NH2	1.82	0.74
17:AR:46:LYS:HB2	17:AR:58:VAL:HG13	1.69	0.74
1:AA:1757:C:OP1	53:B5:2264:U:O5'	2.05	0.74
1:AA:1501:A:C4	1:AA:1546:G:C2	2.74	0.74
17:AR:41:THR:HG23	17:AR:61:PHE:O	1.88	0.74
1:AA:1756:U:H2'	53:B5:2263:C:OP1	1.87	0.74
1:AA:1665:A:P	53:B5:1936:A:H5'	2.27	0.74
17:AR:19:TRP:HB3	17:AR:38:ARG:HD2	1.70	0.74
1:AA:701:U:O4	40:BP:172:ALA:HB3	1.33	0.74
17:AR:44:SER:O	17:AR:58:VAL:HG22	1.87	0.74
14:AO:123:HIS:CE1	53:B5:846:A:N3	2.56	0.73
37:BM:147:TRP:CZ2	37:BM:153:ALA:CB	2.70	0.73
1:AA:1521:G:C4	18:AT:79:LEU:N	2.55	0.73
1:AA:1521:G:H4'	1:AA:1522:A:OP1	1.88	0.73
17:AR:42:LEU:HD13	17:AR:43:ILE:N	2.03	0.73
17:AR:36:ALA:HB1	17:AR:68:VAL:CG1	2.18	0.73
17:AR:40:LYS:HG2	17:AR:66:HIS:C	2.08	0.73
1:AA:702:G:O6	40:BP:172:ALA:HB2	1.82	0.73
1:AA:1521:G:C5'	1:AA:1522:A:OP1	2.34	0.73
17:AR:203:THR:HG23	17:AR:245:PHE:HE1	1.54	0.73
1:AA:1200:A:C6	13:AN:16:LYS:C	2.62	0.73
1:AA:799:A:N3	40:BP:161:ALA:HB3	1.99	0.73
37:BM:147:TRP:HZ2	37:BM:153:ALA:CB	2.02	0.73
1:AA:699:U:C2'	40:BP:170:ALA:N	2.50	0.72
1:AA:972:A:C1'	53:B5:847:A:H2	2.01	0.72
17:AR:262:VAL:HG22	17:AR:272:ASP:O	1.89	0.72
1:AA:700:C:H5	40:BP:169:ALA:O	1.55	0.72
17:AR:199:ILE:HD11	17:AR:201:THR:O	1.90	0.72
17:AR:222:LEU:HD13	17:AR:232:TYR:OH	1.89	0.72
1:AA:1501:A:C4	1:AA:1546:G:N3	2.57	0.72
1:AA:799:A:H4'	40:BP:161:ALA:O	1.87	0.72
1:AA:1653:A:H2	53:B5:2301:U:C2'	1.96	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:B8:43:LYS:CE	35:BK:121:PHE:CE1	2.73	0.72
17:AR:165:ASP:HB2	17:AR:183:LEU:C	2.10	0.72
23:B8:43:LYS:CB	35:BK:121:PHE:CE1	2.73	0.72
17:AR:172:ALA:HB2	17:AR:202:LEU:HG	1.70	0.71
17:AR:34:LEU:HD11	17:AR:80:ALA:HB1	1.71	0.71
17:AR:34:LEU:CD2	17:AR:42:LEU:HD11	2.19	0.71
17:AR:172:ALA:HB1	17:AR:199:ILE:HD12	1.73	0.71
1:AA:698:U:C2	40:BP:164:ALA:C	2.57	0.71
12:AM:110:ARG:NH2	34:BJ:116:TYR:CE1	2.50	0.71
17:AR:275:ARG:C	17:AR:285:ALA:HA	2.11	0.71
17:AR:283:LYS:HB3	17:AR:288:HIS:HA	1.73	0.71
17:AR:164:ASP:OD2	17:AR:168:THR:HG22	1.90	0.71
17:AR:115:ILE:HG22	17:AR:116:ASP:O	1.90	0.71
17:AR:207:ASP:HB3	17:AR:209:THR:HG22	1.72	0.71
1:AA:1461:C:P	19:A7:30:G:H1'	2.31	0.71
12:AM:110:ARG:CZ	34:BJ:116:TYR:CE2	2.71	0.71
1:AA:1657:A:H4'	42:BR:67:PRO:HG2	1.73	0.71
17:AR:66:HIS:CE1	17:AR:67:ILE:HD13	2.26	0.71
1:AA:1633:A:O3'	1:AA:1634:C:H4'	1.91	0.71
1:AA:1756:U:O4'	53:B5:2262:A:C2	2.42	0.71
1:AA:1776:G:O3'	53:B5:2194:G:C5'	2.38	0.70
1:AA:702:G:H8	40:BP:175:ALA:CB	1.93	0.70
1:AA:698:U:C5	40:BP:164:ALA:O	2.43	0.70
1:AA:1573:G:N3	19:A7:41:U:H5''	2.06	0.70
1:AA:1501:A:C1'	1:AA:1546:G:N3	2.53	0.70
37:BM:108:ILE:HG21	37:BM:160:ALA:HB1	1.73	0.70
17:AR:283:LYS:HD3	17:AR:288:HIS:CB	2.22	0.70
17:AR:285:ALA:HB3	17:AR:313:TRP:HZ2	1.54	0.70
17:AR:96:THR:HG23	17:AR:98:GLU:N	2.05	0.70
1:AA:1121:A:H4'	53:B5:2191:U:H5'	1.74	0.70
1:AA:1461:C:OP1	19:A7:30:G:C1'	2.32	0.70
1:AA:702:G:OP1	40:BP:177:ALA:HA	1.91	0.70
1:AA:857:U:H2'	1:AA:858:G:O5'	1.91	0.70
1:AA:1502:G:O6	1:AA:1547:C:C5	2.45	0.70
17:AR:34:LEU:HG	17:AR:42:LEU:CD1	2.14	0.70
17:AR:263:PHE:CE1	17:AR:270:LEU:HG	2.26	0.70
1:AA:1502:G:C8	1:AA:1547:C:O2	2.45	0.69
1:AA:856:A:C2	7:AH:32:LYS:O	2.45	0.69
17:AR:275:ARG:HB3	17:AR:276:PRO:HD2	1.75	0.69
1:AA:1756:U:H4'	53:B5:2262:A:C4	2.28	0.69
1:AA:1521:G:H22	18:AT:78:LYS:CE	2.04	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
17:AR:26:SER:HB2	17:AR:30:PRO:HD2	1.75	0.69
17:AR:34:LEU:HD12	17:AR:73:LEU:HD21	1.75	0.69
17:AR:26:SER:CB	17:AR:30:PRO:HD2	2.23	0.69
17:AR:170:ILE:CG2	17:AR:202:LEU:HD13	2.23	0.69
1:AA:1390:C:OP2	17:AR:281:TYR:OH	2.10	0.69
17:AR:135:THR:HG23	17:AR:137:LYS:H	1.58	0.69
17:AR:164:ASP:H	17:AR:168:THR:N	1.91	0.68
17:AR:203:THR:HG23	17:AR:245:PHE:CE1	2.28	0.68
23:B8:43:LYS:NZ	35:BK:121:PHE:HE1	1.88	0.68
17:AR:173:GLY:H	17:AR:199:ILE:HG23	1.57	0.68
1:AA:702:G:H1'	40:BP:175:ALA:HB1	1.70	0.68
14:AO:149:LEU:CD1	53:B5:846:A:O3'	2.42	0.68
1:AA:1630:C:H5''	1:AA:1636:G:H21	1.51	0.68
17:AR:283:LYS:HE2	17:AR:289:ALA:N	2.09	0.68
32:BH:40:HIS:CG	32:BH:41:ILE:H	2.11	0.68
1:AA:1521:G:C2	18:AT:78:LYS:HD2	2.29	0.68
17:AR:211:ILE:CD1	17:AR:225:LEU:HD13	2.09	0.68
17:AR:314:GLN:HG2	17:AR:315:VAL:N	2.09	0.68
1:AA:699:U:O4'	40:BP:169:ALA:HB3	1.93	0.68
17:AR:240:VAL:HG22	17:AR:256:THR:HG22	1.75	0.68
17:AR:276:PRO:CA	17:AR:285:ALA:HA	2.24	0.68
17:AR:87:LYS:HD2	17:AR:106:HIS:O	1.93	0.68
17:AR:170:ILE:HG22	17:AR:202:LEU:CD1	2.24	0.67
17:AR:34:LEU:HD21	17:AR:71:CYS:SG	2.33	0.67
17:AR:257:ALA:O	17:AR:283:LYS:HA	1.95	0.67
17:AR:169:ILE:CG1	17:AR:183:LEU:HD21	2.25	0.67
1:AA:1521:G:C4'	1:AA:1522:A:OP1	2.37	0.67
17:AR:33:LEU:HD23	17:AR:33:LEU:H	1.59	0.67
37:BM:107:GLY:HA3	37:BM:156:ALA:HB1	1.77	0.67
1:AA:1516:C:C2'	1:AA:1517:U:O5'	2.42	0.67
37:BM:168:ALA:O	37:BM:172:ALA:N	2.28	0.67
1:AA:1390:C:P	17:AR:281:TYR:OH	2.50	0.67
17:AR:259:GLY:HA2	17:AR:284:ALA:N	2.08	0.67
17:AR:199:ILE:HD11	17:AR:202:LEU:HD23	1.77	0.67
1:AA:698:U:C6	40:BP:164:ALA:C	2.62	0.66
17:AR:259:GLY:CA	17:AR:284:ALA:HA	2.22	0.66
17:AR:10:ARG:HG2	17:AR:54:PHE:CE1	2.30	0.66
17:AR:32:LEU:HD22	17:AR:32:LEU:O	1.95	0.66
37:BM:150:ALA:O	37:BM:154:ALA:HB3	1.96	0.66
17:AR:21:THR:HA	17:AR:290:VAL:HG11	1.78	0.66
17:AR:285:ALA:O	17:AR:305:TYR:HE2	1.78	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
17:AR:12:THR:HG22	17:AR:311:ARG:CG	2.23	0.66
17:AR:286:GLU:CD	17:AR:289:ALA:HB2	2.15	0.66
17:AR:203:THR:CG2	17:AR:212:ALA:HB3	2.25	0.66
1:AA:1756:U:H4'	53:B5:2263:C:H1'	1.77	0.66
1:AA:1502:G:N3	1:AA:1547:C:N4	2.44	0.66
17:AR:67:ILE:O	17:AR:84:SER:HB2	1.95	0.65
17:AR:83:ALA:HB1	17:AR:110:VAL:HG12	1.75	0.65
1:AA:921:G:C1'	26:BB:141:PRO:HG3	2.26	0.65
17:AR:93:ASP:CB	17:AR:96:THR:HG22	2.23	0.65
12:AM:110:ARG:NE	34:BJ:116:TYR:CG	2.60	0.65
17:AR:11:GLY:N	17:AR:312:VAL:HG23	2.11	0.65
17:AR:32:LEU:N	17:AR:32:LEU:HD13	2.11	0.65
17:AR:46:LYS:HB2	17:AR:58:VAL:HG11	1.76	0.65
17:AR:81:LEU:HD13	17:AR:91:LEU:HD23	1.78	0.65
23:B8:43:LYS:CE	35:BK:121:PHE:HD1	2.01	0.65
1:AA:629:U:H5'	53:B5:846:A:H1'	1.78	0.65
17:AR:260:ILE:CG1	17:AR:284:ALA:HB1	2.23	0.65
1:AA:1521:G:N7	18:AT:79:LEU:N	2.44	0.65
17:AR:209:THR:CG2	17:AR:210:LEU:HD22	2.10	0.65
17:AR:35:SER:O	17:AR:42:LEU:HD22	1.97	0.65
17:AR:196:ASN:HD21	17:AR:217:ASP:HB2	1.61	0.65
17:AR:283:LYS:HG2	17:AR:284:ALA:H	1.61	0.65
17:AR:38:ARG:O	17:AR:67:ILE:HG13	1.96	0.65
12:AM:110:ARG:CZ	34:BJ:116:TYR:CG	2.80	0.65
1:AA:799:A:H4'	40:BP:161:ALA:C	2.16	0.64
17:AR:33:LEU:HD22	17:AR:47:LEU:HD21	1.78	0.64
19:A7:19:G:N7	53:B5:2454:G:C8	2.65	0.64
1:AA:702:G:P	40:BP:177:ALA:HA	2.38	0.64
1:AA:1756:U:H1'	53:B5:2263:C:OP1	1.96	0.64
17:AR:34:LEU:HD21	17:AR:42:LEU:CD2	2.16	0.64
17:AR:74:THR:HG22	17:AR:79:TYR:H	1.61	0.64
1:AA:1502:G:C6	1:AA:1547:C:N4	2.65	0.64
1:AA:1736:U:C5'	53:B5:1932:A:H61	2.11	0.64
17:AR:7:LEU:HA	17:AR:315:VAL:CG2	2.28	0.64
17:AR:7:LEU:HA	17:AR:315:VAL:HG22	1.78	0.64
1:AA:1460:G:H4'	19:A7:30:G:C5'	2.26	0.64
1:AA:921:G:H5''	26:BB:140:ASN:HD21	1.62	0.64
17:AR:256:THR:O	17:AR:283:LYS:HE3	1.97	0.64
17:AR:90:ARG:NH1	17:AR:99:THR:HG21	2.12	0.64
17:AR:22:SER:O	17:AR:23:LEU:HD23	1.98	0.64
1:AA:1756:U:C2'	53:B5:2263:C:C5'	2.43	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
17:AR:34:LEU:HD11	17:AR:80:ALA:CB	2.28	0.64
1:AA:1755:G:HO2'	53:B5:2262:A:H2	0.65	0.63
17:AR:262:VAL:HG23	17:AR:272:ASP:H	1.63	0.63
1:AA:1630:C:H5'	1:AA:1636:G:N2	2.08	0.63
17:AR:274:LEU:HD13	17:AR:313:TRP:CD2	2.33	0.63
17:AR:286:GLU:HA	17:AR:305:TYR:HD2	1.63	0.63
1:AA:1501:A:O4'	1:AA:1546:G:N2	2.31	0.63
1:AA:1636:G:O6	1:AA:1762:C:O2	2.16	0.63
17:AR:13:LEU:N	17:AR:13:LEU:HD22	2.13	0.63
17:AR:274:LEU:O	17:AR:284:ALA:HA	1.99	0.63
17:AR:40:LYS:HG2	17:AR:66:HIS:O	1.96	0.63
1:AA:1756:U:C4'	53:B5:2262:A:C4	2.81	0.63
17:AR:214:ALA:HB1	17:AR:240:VAL:HB	1.81	0.63
1:AA:701:U:O4	40:BP:169:ALA:C	2.35	0.63
37:BM:168:ALA:O	37:BM:172:ALA:HB2	1.99	0.63
1:AA:972:A:H5'	53:B5:848:A:H2	1.57	0.63
1:AA:1516:C:H2'	1:AA:1517:U:O5'	1.99	0.63
17:AR:164:ASP:C	17:AR:183:LEU:HB2	2.19	0.63
1:AA:1736:U:H5'	53:B5:1932:A:N6	2.13	0.63
1:AA:1776:G:C3'	53:B5:2194:G:H5''	2.26	0.62
17:AR:181:TRP:CZ3	17:AR:188:ILE:HB	2.34	0.62
17:AR:86:ASP:O	17:AR:87:LYS:HG2	1.99	0.62
1:AA:699:U:O2	40:BP:167:ALA:O	2.17	0.62
1:AA:699:U:N3	40:BP:167:ALA:C	2.50	0.62
1:AA:701:U:O4	40:BP:169:ALA:CA	2.47	0.62
17:AR:19:TRP:CZ3	17:AR:306:THR:HG22	2.34	0.62
17:AR:170:ILE:N	17:AR:170:ILE:HD12	2.14	0.62
1:AA:1120:C:O2'	53:B5:2190:U:H5''	1.99	0.62
37:BM:168:ALA:CA	37:BM:171:ALA:HB3	2.29	0.62
1:AA:972:A:C4'	53:B5:847:A:C2	2.83	0.62
17:AR:238:ASP:OD2	17:AR:258:THR:HB	2.00	0.62
1:AA:1460:G:O3'	19:A7:30:G:C4'	2.47	0.62
17:AR:178:VAL:HB	17:AR:192:PHE:HB2	1.81	0.62
1:AA:858:G:H3'	7:AH:27:ILE:HB	1.79	0.61
17:AR:86:ASP:CG	17:AR:88:THR:HG22	2.19	0.61
17:AR:175:ASP:O	17:AR:176:LYS:HG2	2.00	0.61
1:AA:1574:A:C2	19:A7:41:U:C1'	2.83	0.61
1:AA:921:G:H5''	26:BB:140:ASN:ND2	2.15	0.61
17:AR:286:GLU:HG2	17:AR:289:ALA:N	2.15	0.61
1:AA:1519:G:C4'	1:AA:1520:U:OP1	2.44	0.61
17:AR:283:LYS:HZ3	17:AR:288:HIS:CD2	2.18	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
17:AR:131:ILE:CD1	17:AR:151:VAL:HG11	2.31	0.61
17:AR:242:SER:OG	17:AR:292:LEU:HD22	2.01	0.61
17:AR:100:TYR:CD2	17:AR:101:GLN:HG3	2.36	0.61
1:AA:1653:A:C2	53:B5:2301:U:C2'	2.82	0.61
17:AR:256:THR:O	17:AR:283:LYS:HG3	2.01	0.61
17:AR:288:HIS:HB3	17:AR:306:THR:HG21	1.82	0.61
17:AR:287:PRO:HG3	17:AR:307:ASP:HB3	1.83	0.61
17:AR:9:LEU:HD21	17:AR:311:ARG:HB3	1.83	0.61
1:AA:1121:A:H1'	53:B5:2191:U:OP1	2.01	0.61
17:AR:220:ILE:N	17:AR:220:ILE:HD12	2.16	0.61
37:BM:111:PRO:HA	37:BM:163:ALA:HB1	0.62	0.61
37:BM:111:PRO:N	37:BM:163:ALA:HB1	2.04	0.61
1:AA:699:U:C4'	40:BP:166:ALA:HA	2.31	0.61
17:AR:276:PRO:HB2	17:AR:278:PHE:CE1	2.35	0.61
1:AA:1665:A:OP1	53:B5:1935:G:O2'	2.19	0.61
23:B8:43:LYS:CD	35:BK:121:PHE:HE1	2.12	0.61
17:AR:8:VAL:HG12	17:AR:9:LEU:N	2.16	0.60
17:AR:164:ASP:HB2	17:AR:168:THR:N	2.16	0.60
14:AO:149:LEU:CD1	53:B5:847:A:H5'	2.14	0.60
17:AR:238:ASP:CG	17:AR:258:THR:HB	2.20	0.60
17:AR:144:LEU:N	17:AR:144:LEU:HD12	2.16	0.60
17:AR:30:PRO:HB2	17:AR:32:LEU:HD11	1.82	0.60
17:AR:222:LEU:HD23	17:AR:223:TRP:N	2.16	0.60
17:AR:223:TRP:CZ3	17:AR:230:ALA:HB2	2.37	0.60
17:AR:156:VAL:HB	17:AR:167:VAL:HG21	1.80	0.60
1:AA:1200:A:C2	13:AN:16:LYS:CB	2.84	0.60
1:AA:701:U:H2'	40:BP:176:ALA:O	2.02	0.60
17:AR:262:VAL:CG2	17:AR:271:VAL:HG12	2.31	0.60
37:BM:111:PRO:HB3	37:BM:163:ALA:O	2.01	0.59
17:AR:164:ASP:H	17:AR:167:VAL:C	2.05	0.59
17:AR:266:ASP:CB	17:AR:267:PRO:HD3	2.26	0.59
1:AA:1501:A:C8	1:AA:1546:G:C4	2.90	0.59
17:AR:41:THR:HG22	17:AR:42:LEU:N	2.17	0.59
17:AR:193:ILE:H	17:AR:193:ILE:CD1	2.12	0.59
12:AM:110:ARG:CZ	34:BJ:116:TYR:CD2	2.86	0.59
17:AR:264:SER:HB3	17:AR:267:PRO:HD2	1.83	0.59
1:AA:1579:C:H3'	8:AI:131:GLY:H	1.67	0.59
17:AR:11:GLY:O	17:AR:312:VAL:HG22	2.01	0.59
17:AR:33:LEU:HD23	17:AR:33:LEU:N	2.17	0.59
1:AA:1778:G:O5'	53:B5:2274:U:C5'	2.47	0.59
37:BM:111:PRO:CG	37:BM:163:ALA:HA	2.33	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:699:U:C5'	40:BP:166:ALA:HA	2.32	0.59
1:AA:703:G:OP2	40:BP:176:ALA:HB1	2.01	0.59
1:AA:1521:G:H21	18:AT:78:LYS:NZ	2.00	0.59
17:AR:42:LEU:HD13	17:AR:43:ILE:H	1.68	0.59
17:AR:201:THR:O	17:AR:202:LEU:HD23	2.02	0.59
1:AA:864:U:H5''	1:AA:865:A:H5'	1.85	0.59
1:AA:995:U:OP1	53:B5:2197:C:OP2	2.21	0.59
53:B5:1018:G:H1	53:B5:1034:U:H3	1.51	0.59
17:AR:284:ALA:HB3	17:AR:286:GLU:OE1	2.01	0.59
17:AR:199:ILE:HD11	17:AR:202:LEU:CD2	2.32	0.59
1:AA:971:G:H21	53:B5:846:A:H62	1.43	0.59
1:AA:1200:A:C6	13:AN:16:LYS:HA	2.37	0.58
1:AA:856:A:C6	7:AH:32:LYS:CA	2.85	0.58
1:AA:1775:G:H5''	53:B5:2193:U:OP1	2.03	0.58
1:AA:1573:G:N2	19:A7:42:G:P	2.57	0.58
1:AA:1574:A:C5	19:A7:41:U:H4'	2.23	0.58
1:AA:1777:U:P	53:B5:2194:G:C4'	2.82	0.58
1:AA:1746:G:P	53:B5:2303:A:H3'	2.43	0.58
1:AA:1637:C:C2'	1:AA:1638:C:O5'	2.49	0.58
1:AA:1502:G:N9	1:AA:1547:C:N3	2.51	0.58
1:AA:1744:A:C2	53:B5:2302:G:H8	2.21	0.58
17:AR:72:THR:HG22	17:AR:73:LEU:N	2.18	0.58
17:AR:197:SER:HB2	17:AR:216:LYS:HB3	1.83	0.58
37:BM:148:LYS:HB2	37:BM:149:ALA:HB3	1.85	0.58
17:AR:42:LEU:HD21	17:AR:71:CYS:CB	2.34	0.58
1:AA:1637:C:H2'	1:AA:1638:C:O5'	2.04	0.58
1:AA:972:A:P	53:B5:848:A:H2	2.27	0.58
17:AR:106:HIS:CE1	17:AR:132:LYS:HD2	2.39	0.58
17:AR:21:THR:HG22	17:AR:36:ALA:C	2.23	0.58
17:AR:114:ASP:OD1	17:AR:154:VAL:HG23	2.03	0.58
17:AR:224:ASN:ND2	17:AR:227:ALA:HB3	2.19	0.58
1:AA:698:U:N3	40:BP:164:ALA:O	2.35	0.57
17:AR:33:LEU:HD21	17:AR:45:TRP:HB2	1.86	0.57
1:AA:1573:G:N2	19:A7:41:U:O5'	2.37	0.57
17:AR:283:LYS:HG2	17:AR:289:ALA:H	1.69	0.57
17:AR:135:THR:HG23	17:AR:137:LYS:N	2.18	0.57
1:AA:1390:C:P	17:AR:281:TYR:HH	2.28	0.57
17:AR:166:SER:O	17:AR:167:VAL:HG13	2.05	0.57
17:AR:169:ILE:HD11	17:AR:181:TRP:CD1	2.40	0.57
1:AA:1501:A:N6	1:AA:1546:G:H8	1.97	0.57
17:AR:301:LEU:HB3	17:AR:313:TRP:HB2	1.85	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:702:G:C4	40:BP:176:ALA:HB3	2.32	0.57
17:AR:283:LYS:HE2	17:AR:289:ALA:O	2.04	0.57
17:AR:164:ASP:H	17:AR:167:VAL:CA	2.17	0.57
17:AR:22:SER:HA	17:AR:291:SER:OG	2.04	0.57
27:BC:99:LEU:HD21	27:BC:159:ARG:HH22	1.68	0.57
1:AA:702:G:C4	40:BP:176:ALA:CB	2.88	0.57
1:AA:1521:G:O5'	1:AA:1522:A:OP2	2.23	0.57
17:AR:21:THR:HG22	17:AR:36:ALA:O	2.05	0.57
17:AR:262:VAL:HG23	17:AR:271:VAL:CG1	2.34	0.57
17:AR:34:LEU:HD11	17:AR:71:CYS:SG	2.45	0.57
37:BM:150:ALA:O	37:BM:154:ALA:CB	2.53	0.57
37:BM:148:LYS:HB2	37:BM:149:ALA:HB2	1.84	0.57
1:AA:701:U:O2	40:BP:171:ALA:O	2.22	0.57
17:AR:286:GLU:CG	17:AR:305:TYR:CD2	2.87	0.57
17:AR:31:ASN:HA	17:AR:47:LEU:HD12	1.87	0.56
1:AA:1756:U:C2'	53:B5:2263:C:P	2.94	0.56
17:AR:106:HIS:HA	17:AR:132:LYS:HE3	1.87	0.56
1:AA:799:A:H5'	40:BP:165:ALA:N	2.20	0.56
17:AR:10:ARG:HB3	17:AR:312:VAL:HG23	1.88	0.56
17:AR:34:LEU:HD23	17:AR:35:SER:N	2.20	0.56
17:AR:156:VAL:HB	17:AR:167:VAL:HG23	1.85	0.56
17:AR:83:ALA:CB	17:AR:113:VAL:HG13	2.25	0.56
17:AR:141:LEU:N	17:AR:141:LEU:HD12	2.21	0.56
1:AA:972:A:C4'	53:B5:847:A:H2	2.18	0.56
17:AR:224:ASN:CG	17:AR:227:ALA:HB3	2.26	0.56
20:B0:38:ILE:HG23	20:B0:39:ASP:H	1.69	0.56
17:AR:284:ALA:HB3	17:AR:289:ALA:HB2	1.86	0.56
17:AR:163:ASP:HA	17:AR:167:VAL:HA	1.87	0.56
1:AA:858:G:O4'	1:AA:858:G:OP2	2.24	0.56
17:AR:237:GLN:HG2	17:AR:237:GLN:O	2.06	0.56
1:AA:1777:U:H5''	53:B5:2194:G:O2'	2.05	0.56
1:AA:1756:U:H4'	53:B5:2262:A:C5	2.41	0.56
1:AA:1776:G:C3'	53:B5:2194:G:C5'	2.82	0.56
17:AR:259:GLY:HA2	17:AR:283:LYS:C	2.25	0.56
26:BB:50:HIS:H	53:B5:1796:G:H5''	1.71	0.56
1:AA:1777:U:C5'	53:B5:2194:G:O2'	2.54	0.56
17:AR:275:ARG:HB3	17:AR:276:PRO:CD	2.35	0.56
17:AR:305:TYR:HB2	17:AR:309:VAL:O	2.06	0.56
37:BM:106:GLU:OE1	37:BM:153:ALA:O	2.24	0.56
1:AA:867:G:H5''	1:AA:868:G:OP2	2.06	0.56
1:AA:1461:C:P	19:A7:30:G:C1'	2.94	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
14:AO:149:LEU:HD12	53:B5:846:A:O3'	2.04	0.55
1:AA:1518:U:H5''	1:AA:1519:G:OP2	2.06	0.55
17:AR:179:LYS:HD3	17:AR:181:TRP:CZ2	2.41	0.55
17:AR:9:LEU:CD2	17:AR:311:ARG:HB3	2.36	0.55
17:AR:34:LEU:HD12	17:AR:73:LEU:CD2	2.35	0.55
7:AH:32:LYS:C	7:AH:33:VAL:N	2.59	0.55
1:AA:1631:A:OP1	1:AA:1636:G:O2'	2.13	0.55
17:AR:302:PHE:CD2	17:AR:312:VAL:HG12	2.42	0.55
1:AA:858:G:C4	7:AH:61:ILE:CG1	2.88	0.55
6:AG:72:HIS:CD2	6:AG:89:ILE:H	2.24	0.55
1:AA:972:A:H5''	53:B5:848:A:N3	2.14	0.55
17:AR:276:PRO:HB2	17:AR:278:PHE:CD1	2.40	0.55
17:AR:169:ILE:HD12	17:AR:169:ILE:C	2.27	0.55
18:AT:28:LEU:HD21	18:AT:111:ILE:HD11	1.88	0.55
17:AR:34:LEU:C	17:AR:34:LEU:HD23	2.27	0.55
1:AA:1734:G:H21	53:B5:1934:G:C4'	2.06	0.55
17:AR:283:LYS:CG	17:AR:284:ALA:H	2.19	0.55
17:AR:210:LEU:HD12	17:AR:222:LEU:CD2	2.26	0.55
23:B8:43:LYS:HB2	35:BK:121:PHE:CE1	2.41	0.55
1:AA:1756:U:H1'	53:B5:2262:A:O2'	2.05	0.55
17:AR:264:SER:CB	17:AR:267:PRO:HD2	2.37	0.55
1:AA:699:U:C2'	40:BP:169:ALA:C	2.72	0.55
17:AR:315:VAL:HG12	17:AR:316:MET:N	2.22	0.55
17:AR:19:TRP:CG	17:AR:38:ARG:HD2	2.42	0.55
17:AR:199:ILE:C	17:AR:199:ILE:HD13	2.28	0.54
1:AA:1521:G:N2	18:AT:78:LYS:HZ3	2.03	0.54
17:AR:250:TYR:O	17:AR:265:LEU:HG	2.06	0.54
17:AR:222:LEU:C	17:AR:222:LEU:HD23	2.27	0.54
37:BM:155:ALA:O	37:BM:157:ALA:N	2.41	0.54
19:A7:56:C:N4	53:B5:2484:A:N3	2.55	0.54
1:AA:1756:U:C5'	53:B5:2262:A:C6	2.88	0.54
17:AR:31:ASN:C	17:AR:32:LEU:HD13	2.28	0.54
17:AR:150:TRP:HB2	17:AR:174:ASN:ND2	2.22	0.54
1:AA:702:G:C5	40:BP:172:ALA:HA	2.23	0.54
1:AA:1501:A:C8	1:AA:1546:G:N3	2.76	0.54
1:AA:1521:G:C8	18:AT:79:LEU:N	2.75	0.54
17:AR:21:THR:HG23	17:AR:36:ALA:HB3	1.90	0.54
37:BM:111:PRO:CB	37:BM:163:ALA:C	2.68	0.54
17:AR:188:ILE:HG23	17:AR:188:ILE:O	2.07	0.54
1:AA:1200:A:C2	13:AN:16:LYS:HB2	2.42	0.54
1:AA:1735:G:C5'	53:B5:1933:A:O2'	2.56	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
17:AR:276:PRO:HA	17:AR:285:ALA:CA	2.38	0.54
1:AA:799:A:O3'	40:BP:164:ALA:HB2	2.08	0.54
17:AR:262:VAL:HG23	17:AR:272:ASP:N	2.23	0.54
17:AR:199:ILE:HG23	17:AR:199:ILE:O	2.08	0.53
17:AR:248:ASN:ND2	17:AR:249:ARG:HG2	2.22	0.53
1:AA:858:G:C3'	7:AH:27:ILE:HB	2.38	0.53
1:AA:864:U:C5'	1:AA:865:A:H5'	2.37	0.53
1:AA:701:U:O2	40:BP:171:ALA:C	2.45	0.53
17:AR:13:LEU:HB2	17:AR:310:ILE:HB	1.89	0.53
1:AA:1664:U:H1'	53:B5:1934:G:O2'	2.09	0.53
17:AR:270:LEU:HD12	17:AR:270:LEU:N	2.24	0.53
37:BM:111:PRO:CG	37:BM:163:ALA:CB	2.87	0.53
1:AA:1521:G:H21	18:AT:78:LYS:HZ3	1.56	0.53
1:AA:701:U:C4'	40:BP:173:ALA:HB1	2.05	0.53
17:AR:263:PHE:HE1	17:AR:270:LEU:CD1	2.22	0.53
17:AR:193:ILE:O	17:AR:223:TRP:HH2	1.91	0.53
17:AR:136:ILE:HG23	17:AR:137:LYS:N	2.23	0.53
17:AR:220:ILE:HB	17:AR:234:LEU:HB2	1.90	0.53
17:AR:170:ILE:HG22	17:AR:171:SER:N	2.24	0.53
19:A7:19:G:N1	53:B5:2454:G:C6	2.69	0.53
1:AA:1574:A:H2'	19:A7:40:5MC:HO2'	1.68	0.53
17:AR:276:PRO:HB3	17:AR:286:GLU:N	2.23	0.53
17:AR:98:GLU:HG2	17:AR:99:THR:N	2.24	0.53
17:AR:178:VAL:HG13	17:AR:202:LEU:HD11	1.91	0.53
17:AR:164:ASP:H	17:AR:167:VAL:HA	1.74	0.53
1:AA:1339:C:H5'	17:AR:102:ARG:HH21	1.68	0.53
21:B1:40:LYS:H	53:B5:355:A:H4'	1.74	0.53
17:AR:284:ALA:C	17:AR:286:GLU:H	2.13	0.52
1:AA:858:G:C2	7:AH:61:ILE:CA	2.92	0.52
1:AA:1744:A:N3	53:B5:2302:G:C8	2.77	0.52
19:A7:56:C:N4	53:B5:2484:A:C2	2.77	0.52
17:AR:172:ALA:HB1	17:AR:199:ILE:CD1	2.40	0.52
40:BP:134:HIS:CG	40:BP:135:LYS:H	2.27	0.52
1:AA:1755:G:C2'	53:B5:2262:A:C2	2.70	0.52
1:AA:1735:G:H4'	53:B5:1933:A:C2'	2.33	0.52
17:AR:287:PRO:HG2	17:AR:306:THR:OG1	2.09	0.52
17:AR:170:ILE:HG12	17:AR:211:ILE:HD13	1.90	0.52
17:AR:85:TRP:HB3	17:AR:109:ASP:OD1	2.10	0.52
17:AR:263:PHE:HE1	17:AR:270:LEU:HD11	1.75	0.52
1:AA:1502:G:N1	1:AA:1547:C:N4	2.57	0.52
1:AA:1120:C:O2'	53:B5:2190:U:H4'	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:BB:5:ILE:HD13	26:BB:5:ILE:H	1.75	0.52
19:A7:59:U:C5	19:A7:60:C:C4	2.98	0.52
17:AR:74:THR:HG23	17:AR:77:GLY:H	1.75	0.52
17:AR:270:LEU:HD12	17:AR:270:LEU:H	1.75	0.52
1:AA:722:G:H22	1:AA:732:G:H1	1.58	0.52
17:AR:256:THR:OG1	17:AR:258:THR:HG22	2.10	0.52
19:A7:56:C:O4'	53:B5:2484:A:H4'	2.10	0.52
1:AA:1460:G:H5'	19:A7:29:A:O2'	2.10	0.52
17:AR:69:GLN:HG2	17:AR:85:TRP:CD1	2.45	0.52
17:AR:9:LEU:HD13	17:AR:9:LEU:C	2.29	0.52
17:AR:149:ASP:HB3	17:AR:174:ASN:HB2	1.92	0.52
1:AA:1657:A:C5'	42:BR:67:PRO:HD2	2.28	0.51
17:AR:285:ALA:CB	17:AR:313:TRP:CH2	2.93	0.51
1:AA:1777:U:OP1	53:B5:2194:G:O2'	2.22	0.51
17:AR:283:LYS:HD3	17:AR:288:HIS:CD2	2.44	0.51
17:AR:118:LYS:HD2	17:AR:118:LYS:N	2.25	0.51
17:AR:19:TRP:HE1	17:AR:290:VAL:HG21	1.76	0.51
17:AR:93:ASP:O	17:AR:96:THR:HG22	2.10	0.51
1:AA:922:A:H5''	26:BB:109:GLU:OE1	2.06	0.51
37:BM:168:ALA:O	37:BM:172:ALA:CB	2.58	0.51
17:AR:26:SER:HB3	17:AR:30:PRO:HD2	1.92	0.51
17:AR:66:HIS:ND1	17:AR:67:ILE:HD13	2.25	0.51
1:AA:1521:G:C6	18:AT:79:LEU:N	2.76	0.51
1:AA:1744:A:N3	53:B5:2302:G:H8	2.08	0.51
17:AR:89:LEU:HB2	17:AR:103:PHE:HB2	1.93	0.51
1:AA:1636:G:H2'	1:AA:1637:C:O4'	2.11	0.51
17:AR:30:PRO:HB2	17:AR:32:LEU:CD1	2.41	0.51
17:AR:292:LEU:O	17:AR:292:LEU:HD23	2.11	0.51
23:B8:43:LYS:HG2	35:BK:121:PHE:CE1	2.45	0.50
1:AA:1501:A:N7	1:AA:1546:G:N9	2.59	0.50
1:AA:699:U:C2	40:BP:168:ALA:N	2.79	0.50
1:AA:799:A:H4'	40:BP:164:ALA:H	1.76	0.50
1:AA:857:U:C2'	1:AA:858:G:O5'	2.59	0.50
17:AR:157:VAL:O	17:AR:167:VAL:HG21	2.11	0.50
17:AR:173:GLY:N	17:AR:199:ILE:HG23	2.26	0.50
15:AQ:131:ILE:HD13	15:AQ:131:ILE:H	1.74	0.50
17:AR:276:PRO:HA	17:AR:285:ALA:HA	1.93	0.50
17:AR:305:TYR:HD2	17:AR:311:ARG:HH21	1.56	0.50
1:AA:1460:G:C4'	19:A7:30:G:C5'	2.84	0.50
17:AR:290:VAL:HG13	17:AR:290:VAL:O	2.06	0.50
17:AR:114:ASP:OD2	17:AR:156:VAL:HG13	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
17:AR:81:LEU:HD21	17:AR:122:ILE:CG2	2.41	0.50
1:AA:1776:G:OP1	53:B5:2193:U:C4'	2.53	0.50
1:AA:972:A:C1'	53:B5:847:A:C2	2.87	0.50
17:AR:19:TRP:CZ3	17:AR:306:THR:CG2	2.95	0.50
17:AR:285:ALA:HB3	17:AR:313:TRP:CH2	2.47	0.50
17:AR:314:GLN:HG2	17:AR:315:VAL:H	1.76	0.50
17:AR:99:THR:O	17:AR:99:THR:HG23	2.11	0.50
17:AR:176:LYS:HD3	17:AR:195:HIS:O	2.11	0.50
1:AA:971:G:N2	53:B5:846:A:H61	2.05	0.50
17:AR:305:TYR:CD1	17:AR:305:TYR:N	2.80	0.50
17:AR:7:LEU:N	17:AR:7:LEU:HD12	2.27	0.50
37:BM:147:TRP:HZ2	37:BM:153:ALA:HB1	1.73	0.50
17:AR:169:ILE:CD1	17:AR:181:TRP:CD1	2.95	0.50
17:AR:89:LEU:HD21	17:AR:124:SER:CB	2.34	0.50
33:BI:121:LYS:H	33:BI:121:LYS:HD2	1.76	0.50
17:AR:96:THR:CG2	17:AR:98:GLU:H	2.20	0.50
1:AA:701:U:C5'	40:BP:173:ALA:HB1	2.41	0.49
17:AR:256:THR:C	17:AR:283:LYS:HE3	2.32	0.49
17:AR:115:ILE:HG22	17:AR:119:ALA:HA	1.93	0.49
17:AR:164:ASP:HB3	17:AR:182:ASN:C	2.31	0.49
17:AR:81:LEU:HD21	17:AR:122:ILE:HG21	1.93	0.49
1:AA:699:U:C6	40:BP:165:ALA:O	2.66	0.49
1:AA:1744:A:C2	53:B5:2302:G:C5'	2.77	0.49
17:AR:277:GLU:O	17:AR:278:PHE:HB3	2.12	0.49
17:AR:44:SER:OG	17:AR:58:VAL:HG23	2.11	0.49
17:AR:13:LEU:HD23	17:AR:310:ILE:CG2	2.43	0.49
1:AA:1338:A:O3'	17:AR:102:ARG:NH2	2.44	0.49
1:AA:1665:A:OP1	53:B5:1936:A:O5'	2.28	0.49
17:AR:263:PHE:CE1	17:AR:270:LEU:CD1	2.95	0.49
1:AA:700:C:C5	40:BP:173:ALA:HB3	2.47	0.49
17:AR:69:GLN:CG	17:AR:85:TRP:HE1	2.26	0.49
17:AR:29:GLN:OE1	17:AR:297:ASP:HB3	2.12	0.49
17:AR:131:ILE:HD11	17:AR:151:VAL:HG11	1.94	0.49
53:B5:512:U:H3	53:B5:579:G:H1	1.61	0.49
17:AR:185:GLN:O	17:AR:186:PHE:HB2	2.12	0.49
1:AA:972:A:O4'	53:B5:847:A:N1	2.43	0.49
17:AR:159:ASN:O	17:AR:167:VAL:HG12	2.13	0.49
17:AR:52:GLN:HA	17:AR:52:GLN:NE2	2.20	0.49
17:AR:187:GLN:OE1	17:AR:187:GLN:HA	2.13	0.49
17:AR:258:THR:HG23	17:AR:261:LYS:HE2	1.90	0.49
17:AR:286:GLU:CA	17:AR:305:TYR:CD2	2.93	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
17:AR:34:LEU:CD2	17:AR:71:CYS:HB3	2.43	0.49
17:AR:8:VAL:HG12	17:AR:9:LEU:H	1.77	0.49
17:AR:123:ILE:HG22	17:AR:133:VAL:HA	1.95	0.49
17:AR:131:ILE:HD13	17:AR:151:VAL:HG11	1.95	0.49
17:AR:9:LEU:HD13	17:AR:11:GLY:N	2.27	0.49
37:BM:147:TRP:CZ2	37:BM:153:ALA:HB2	2.48	0.49
17:AR:286:GLU:OE2	17:AR:289:ALA:HB2	2.12	0.48
17:AR:29:GLN:O	17:AR:29:GLN:HG2	2.13	0.48
16:AS:69:GLU:HG3	16:AS:70:ASN:H	1.78	0.48
17:AR:178:VAL:HG13	17:AR:202:LEU:CD1	2.43	0.48
17:AR:123:ILE:HD13	17:AR:169:ILE:HG21	1.95	0.48
1:AA:1520:U:C2'	1:AA:1521:G:OP1	2.58	0.48
1:AA:1460:G:O3'	19:A7:30:G:C1'	2.60	0.48
17:AR:54:PHE:CD1	17:AR:312:VAL:HG21	2.49	0.48
17:AR:87:LYS:HD3	17:AR:107:LYS:O	2.13	0.48
1:AA:1552:U:OP1	13:AN:14:PHE:CG	2.67	0.48
17:AR:33:LEU:HD13	17:AR:302:PHE:CD2	2.48	0.48
32:BH:40:HIS:CG	32:BH:41:ILE:N	2.78	0.48
1:AA:698:U:C4	40:BP:164:ALA:O	2.67	0.48
17:AR:183:LEU:N	17:AR:183:LEU:HD22	2.28	0.48
14:AO:149:LEU:HB2	53:B5:847:A:OP1	2.14	0.48
17:AR:218:GLY:O	17:AR:236:ALA:HB3	2.14	0.48
17:AR:309:VAL:HG12	17:AR:310:ILE:N	2.29	0.48
17:AR:41:THR:HG22	17:AR:42:LEU:O	2.13	0.48
17:AR:154:VAL:O	17:AR:154:VAL:HG23	2.12	0.48
17:AR:169:ILE:HG23	17:AR:183:LEU:CD2	2.43	0.48
17:AR:286:GLU:HG2	17:AR:288:HIS:C	2.34	0.48
17:AR:100:TYR:CE2	17:AR:101:GLN:HG3	2.49	0.48
1:AA:1779:A:H2'	1:AA:1780:A:C8	2.49	0.48
1:AA:1756:U:H4'	53:B5:2263:C:C1'	2.44	0.47
1:AA:698:U:C6	1:AA:799:A:OP2	2.67	0.47
1:AA:1775:G:C4'	53:B5:2193:U:OP1	2.62	0.47
1:AA:1520:U:O4'	1:AA:1522:A:N7	2.47	0.47
17:AR:109:ASP:HB2	17:AR:127:ARG:HD2	1.95	0.47
17:AR:83:ALA:HB2	17:AR:113:VAL:HG11	1.93	0.47
22:B2:104:LEU:H	22:B2:104:LEU:HD13	1.79	0.47
1:AA:858:G:N1	7:AH:61:ILE:CA	2.77	0.47
17:AR:146:GLY:C	17:AR:179:LYS:HE2	2.33	0.47
19:A7:62:A:H2'	19:A7:63:C:C6	2.50	0.47
1:AA:1200:A:C6	13:AN:15:GLY:HA2	2.48	0.47
17:AR:210:LEU:CD1	17:AR:222:LEU:HD21	2.28	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:702:G:C8	40:BP:172:ALA:HA	2.42	0.47
17:AR:221:MET:SD	17:AR:223:TRP:CZ2	3.07	0.47
17:AR:81:LEU:HD13	17:AR:91:LEU:CD2	2.45	0.47
1:AA:629:U:H5'	53:B5:846:A:C1'	2.44	0.47
17:AR:48:THR:HG22	17:AR:50:ASP:CG	2.35	0.47
1:AA:856:A:C4	7:AH:33:VAL:CA	2.96	0.47
1:AA:698:U:C6	40:BP:165:ALA:O	2.61	0.47
17:AR:276:PRO:HB3	17:AR:286:GLU:O	2.15	0.47
17:AR:11:GLY:H	17:AR:312:VAL:HG23	1.79	0.47
17:AR:10:ARG:NE	17:AR:54:PHE:HE1	2.13	0.47
17:AR:222:LEU:HD13	17:AR:232:TYR:CZ	2.49	0.47
37:BM:151:ALA:O	37:BM:155:ALA:N	2.46	0.47
17:AR:165:ASP:CB	17:AR:183:LEU:CB	2.93	0.47
37:BM:148:LYS:HB3	37:BM:149:ALA:H	1.28	0.47
1:AA:1582:G:H21	8:AI:134:ALA:HB1	1.79	0.47
17:AR:13:LEU:HD13	17:AR:54:PHE:HB3	1.94	0.47
17:AR:227:ALA:O	17:AR:228:LYS:HG3	2.15	0.47
1:AA:1520:U:C6	1:AA:1520:U:C3'	2.97	0.47
17:AR:34:LEU:HD21	17:AR:71:CYS:CB	2.45	0.47
17:AR:66:HIS:CE1	17:AR:85:TRP:CE3	3.03	0.47
1:AA:1120:C:O2'	53:B5:2190:U:C5'	2.62	0.47
28:BD:148:ILE:H	28:BD:148:ILE:HD13	1.79	0.47
17:AR:278:PHE:CD2	17:AR:279:ALA:N	2.83	0.46
17:AR:278:PHE:HZ	17:AR:287:PRO:CA	2.27	0.46
17:AR:67:ILE:HD12	17:AR:67:ILE:N	2.30	0.46
8:AI:42:GLU:HB3	8:AI:74:HIS:CE1	2.49	0.46
17:AR:281:TYR:H	17:AR:281:TYR:HD1	1.61	0.46
17:AR:283:LYS:CD	17:AR:288:HIS:CA	2.81	0.46
1:AA:1778:G:O5'	53:B5:2274:U:H5''	2.14	0.46
17:AR:164:ASP:OD2	17:AR:182:ASN:HA	2.15	0.46
17:AR:262:VAL:HG23	17:AR:262:VAL:O	2.16	0.46
17:AR:185:GLN:HB3	17:AR:187:GLN:HG2	1.96	0.46
17:AR:93:ASP:HB3	17:AR:96:THR:HG21	1.95	0.46
17:AR:87:LYS:HA	17:AR:110:VAL:HG23	1.97	0.46
19:A7:37:YYG:H31	19:A7:37:YYG:H1'	1.95	0.46
1:AA:1200:A:C4	1:AA:1200:A:C2	3.03	0.46
17:AR:283:LYS:HB3	17:AR:288:HIS:CA	2.43	0.46
17:AR:283:LYS:CD	17:AR:288:HIS:CD2	2.98	0.46
1:AA:1521:G:H22	18:AT:78:LYS:HG3	1.76	0.46
1:AA:1746:G:OP2	53:B5:2303:A:H3'	2.16	0.46
17:AR:19:TRP:HB3	17:AR:38:ARG:CD	2.41	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
17:AR:274:LEU:HD13	17:AR:313:TRP:CE3	2.50	0.46
17:AR:285:ALA:CB	17:AR:313:TRP:CZ2	2.95	0.46
12:AM:110:ARG:CD	34:BJ:116:TYR:HD1	2.23	0.46
1:AA:866:G:OP1	1:AA:866:G:H8	1.98	0.46
24:B9:51:ALA:HB3	24:B9:54:ILE:HB	1.97	0.46
19:A7:72:C:H5''	53:B5:2234:G:O6	2.16	0.46
1:AA:702:G:C8	40:BP:172:ALA:O	2.69	0.46
1:AA:1547:C:H2'	1:AA:1547:C:H6	1.60	0.46
1:AA:1744:A:H3'	53:B5:2303:A:OP1	2.16	0.46
17:AR:218:GLY:HA2	17:AR:238:ASP:O	2.16	0.46
17:AR:48:THR:HG22	17:AR:50:ASP:H	1.81	0.46
17:AR:96:THR:HG23	17:AR:97:GLY:N	2.30	0.46
17:AR:157:VAL:HG12	17:AR:167:VAL:HB	1.97	0.46
19:A7:1:G:C6	19:A7:73:A:C2	3.04	0.46
1:AA:1521:G:HO2'	18:AT:83:ALA:HA	1.70	0.46
17:AR:10:ARG:HE	17:AR:54:PHE:HE1	1.60	0.46
1:AA:858:G:H1'	7:AH:34:ILE:HG21	1.97	0.46
1:AA:1460:G:H4'	19:A7:29:A:O2'	2.16	0.46
17:AR:134:TRP:CA	17:AR:141:LEU:HD13	2.46	0.46
17:AR:153:GLN:HG3	17:AR:154:VAL:N	2.31	0.46
1:AA:506:A:H3'	1:AA:507:U:C5'	2.46	0.46
1:AA:700:C:OP2	40:BP:169:ALA:HB1	2.15	0.45
1:AA:858:G:O3'	7:AH:29:PRO:HD2	2.17	0.45
17:AR:108:SER:OG	17:AR:127:ARG:HB2	2.16	0.45
23:B8:43:LYS:CB	35:BK:121:PHE:CD1	2.93	0.45
1:AA:701:U:O2'	40:BP:176:ALA:O	2.34	0.45
17:AR:276:PRO:CA	17:AR:285:ALA:CA	2.94	0.45
17:AR:276:PRO:HG3	17:AR:282:SER:HA	1.97	0.45
17:AR:286:GLU:HG3	17:AR:305:TYR:HA	1.97	0.45
17:AR:69:GLN:CG	17:AR:85:TRP:NE1	2.80	0.45
1:AA:1121:A:C1'	53:B5:2191:U:OP1	2.64	0.45
17:AR:156:VAL:HG12	17:AR:169:ILE:HG22	1.97	0.45
15:AQ:82:ARG:HB3	15:AQ:110:HIS:CE1	2.51	0.45
1:AA:1756:U:O2'	53:B5:2262:A:H3'	2.17	0.45
12:AM:119:ILE:HG22	16:AS:106:GLU:H	1.81	0.45
1:AA:701:U:C4	40:BP:169:ALA:C	2.71	0.45
17:AR:283:LYS:CG	17:AR:284:ALA:N	2.78	0.45
37:BM:155:ALA:O	37:BM:156:ALA:C	2.54	0.45
12:AM:120:ARG:HD3	16:AS:103:ASN:H	1.80	0.45
17:AR:284:ALA:CB	17:AR:289:ALA:HB2	2.46	0.45
46:BV:113:GLY:HA3	53:B5:716:G:C4	2.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:856:A:C6	7:AH:32:LYS:N	2.85	0.45
1:AA:1630:C:H5'	1:AA:1636:G:H22	1.82	0.45
17:AR:10:ARG:CG	17:AR:54:PHE:CE1	2.97	0.45
17:AR:31:ASN:HB3	17:AR:47:LEU:HB2	1.99	0.45
53:B5:1003:A:H61	53:B5:1046:A:H61	1.64	0.45
1:AA:1200:A:N3	13:AN:16:LYS:CB	2.79	0.45
37:BM:37:ARG:NH2	37:BM:157:ALA:HA	2.32	0.45
17:AR:169:ILE:O	17:AR:169:ILE:HD12	2.17	0.45
17:AR:262:VAL:CG2	17:AR:271:VAL:CG1	2.94	0.45
26:BB:30:ARG:HG2	26:BB:31:THR:H	1.81	0.45
1:AA:1584:A:C8	8:AI:127:LYS:HB2	2.52	0.45
37:BM:111:PRO:CB	37:BM:163:ALA:O	2.65	0.45
17:AR:164:ASP:N	17:AR:167:VAL:HA	2.32	0.45
26:BB:77:ILE:HD11	26:BB:169:ILE:HD11	1.98	0.45
1:AA:856:A:C2	7:AH:33:VAL:HB	2.52	0.45
1:AA:856:A:N3	7:AH:33:VAL:CA	2.79	0.45
1:AA:1521:G:H21	18:AT:78:LYS:HD2	1.69	0.45
3:AC:164:VAL:HG13	3:AC:165:ASN:H	1.82	0.44
17:AR:34:LEU:CD1	17:AR:73:LEU:CD2	2.95	0.44
17:AR:34:LEU:CD1	17:AR:80:ALA:HB2	2.47	0.44
1:AA:1777:U:OP1	53:B5:2194:G:C2'	2.65	0.44
17:AR:106:HIS:ND1	17:AR:126:SER:HB2	2.32	0.44
1:AA:972:A:H4'	53:B5:847:A:C2	2.52	0.44
17:AR:123:ILE:HA	17:AR:132:LYS:O	2.18	0.44
19:A7:73:A:OP1	53:B5:2603:G:OP1	2.36	0.44
30:BF:118:LYS:HD3	30:BF:118:LYS:H	1.82	0.44
35:BK:121:PHE:CD2	35:BK:121:PHE:CB	2.86	0.44
1:AA:1756:U:C3'	53:B5:2263:C:C1'	2.89	0.44
17:AR:259:GLY:HA2	17:AR:283:LYS:O	2.18	0.44
17:AR:278:PHE:CE2	17:AR:280:GLY:N	2.85	0.44
17:AR:133:VAL:CG1	17:AR:141:LEU:HB2	2.40	0.44
17:AR:91:LEU:HG	17:AR:103:PHE:CE1	2.52	0.44
18:AT:22:LEU:HD12	18:AT:55:TYR:HA	2.00	0.44
17:AR:286:GLU:CD	17:AR:305:TYR:CZ	2.91	0.44
17:AR:172:ALA:HB1	17:AR:199:ILE:O	2.18	0.44
37:BM:108:ILE:HG23	37:BM:160:ALA:CA	2.39	0.44
17:AR:68:VAL:HA	17:AR:84:SER:HB3	2.00	0.44
17:AR:8:VAL:CG1	17:AR:9:LEU:N	2.80	0.44
17:AR:115:ILE:HG22	17:AR:116:ASP:N	2.32	0.44
19:A7:70:C:O2'	53:B5:2236:G:H4'	2.18	0.44
32:BH:129:ARG:H	32:BH:157:ASN:HD21	1.66	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
17:AR:38:ARG:CA	17:AR:67:ILE:HG23	2.22	0.44
17:AR:34:LEU:CD1	17:AR:80:ALA:CB	2.94	0.44
1:AA:1778:G:OP1	53:B5:2274:U:C5'	2.34	0.44
17:AR:56:VAL:HB	17:AR:57:PRO:HD2	1.99	0.44
17:AR:42:LEU:HD21	17:AR:71:CYS:SG	2.58	0.44
17:AR:91:LEU:HG	17:AR:103:PHE:HE1	1.83	0.44
1:AA:1516:C:H6	1:AA:1516:C:O5'	2.01	0.44
1:AA:1501:A:C4'	1:AA:1546:G:N2	2.81	0.43
1:AA:856:A:N3	7:AH:32:LYS:O	2.50	0.43
17:AR:54:PHE:CE2	17:AR:302:PHE:CE2	3.06	0.43
1:AA:1776:G:P	53:B5:2193:U:C5'	2.94	0.43
17:AR:68:VAL:HG23	17:AR:84:SER:HB3	2.00	0.43
1:AA:1520:U:H1'	1:AA:1522:A:H62	1.82	0.43
17:AR:47:LEU:HD22	17:AR:54:PHE:CE2	2.53	0.43
17:AR:262:VAL:CG2	17:AR:272:ASP:N	2.81	0.43
1:AA:1581:A:C8	8:AI:137:ARG:HA	2.54	0.43
17:AR:161:LYS:HA	17:AR:161:LYS:HE2	2.00	0.43
27:BC:342:LEU:H	27:BC:342:LEU:HD23	1.83	0.43
17:AR:193:ILE:N	17:AR:193:ILE:HD13	2.20	0.43
37:BM:155:ALA:C	37:BM:157:ALA:N	2.72	0.43
17:AR:116:ASP:OD2	17:AR:166:SER:HB2	2.19	0.43
17:AR:196:ASN:OD1	17:AR:217:ASP:HB3	2.19	0.43
17:AR:41:THR:OG1	17:AR:62:LYS:HG2	2.18	0.43
42:BR:22:ILE:H	53:B5:1899:G:H5''	1.83	0.43
1:AA:699:U:N3	40:BP:171:ALA:N	2.57	0.43
1:AA:857:U:C6	7:AH:33:VAL:O	2.71	0.43
17:AR:48:THR:H	17:AR:55:GLY:HA2	1.84	0.43
17:AR:74:THR:HG23	17:AR:76:ASP:N	2.34	0.43
1:AA:858:G:C4	7:AH:61:ILE:HG12	2.53	0.43
17:AR:13:LEU:CD1	17:AR:55:GLY:N	2.82	0.43
17:AR:285:ALA:CB	17:AR:313:TRP:HH2	2.32	0.43
17:AR:300:THR:HG23	17:AR:313:TRP:O	2.19	0.43
1:AA:700:C:C6	40:BP:173:ALA:CB	2.98	0.43
1:AA:1777:U:OP1	53:B5:2194:G:C4'	2.66	0.43
33:BI:140:THR:HG22	33:BI:141:LYS:H	1.83	0.43
1:AA:1520:U:H4'	1:AA:1522:A:H8	1.75	0.43
17:AR:19:TRP:NE1	17:AR:290:VAL:HG21	2.34	0.43
17:AR:54:PHE:CE2	17:AR:302:PHE:HE2	2.36	0.43
37:BM:106:GLU:OE1	37:BM:156:ALA:CB	2.67	0.43
1:AA:1734:G:C2	53:B5:1934:G:H4'	2.50	0.43
17:AR:240:VAL:HG22	17:AR:256:THR:CG2	2.46	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
17:AR:133:VAL:HG12	17:AR:142:ALA:N	2.33	0.43
1:AA:796:A:C2'	1:AA:797:G:O5'	2.67	0.43
1:AA:1755:G:H2'	53:B5:2262:A:H2	1.77	0.42
17:AR:10:ARG:HG2	17:AR:54:PHE:HE1	1.81	0.42
17:AR:71:CYS:SG	17:AR:80:ALA:HB1	2.59	0.42
17:AR:172:ALA:HB2	17:AR:199:ILE:HD12	1.96	0.42
17:AR:83:ALA:HB1	17:AR:110:VAL:HG11	1.95	0.42
17:AR:89:LEU:HD12	17:AR:89:LEU:N	2.33	0.42
17:AR:262:VAL:HG21	17:AR:272:ASP:HB3	2.01	0.42
17:AR:292:LEU:CD2	17:AR:292:LEU:N	2.82	0.42
17:AR:185:GLN:NE2	17:AR:187:GLN:HG3	2.34	0.42
42:BR:7:GLN:N	53:B5:3016:A:HO2'	2.17	0.42
12:AM:39:GLY:H	12:AM:42:TYR:HB2	1.85	0.42
1:AA:701:U:O5'	40:BP:173:ALA:HB1	2.18	0.42
1:AA:1664:U:O3'	53:B5:1935:G:C2'	2.62	0.42
17:AR:41:THR:CG2	17:AR:42:LEU:N	2.81	0.42
17:AR:35:SER:CB	17:AR:45:TRP:HE1	2.33	0.42
17:AR:134:TRP:C	17:AR:141:LEU:HD13	2.39	0.42
17:AR:263:PHE:CE1	17:AR:270:LEU:CG	2.99	0.42
17:AR:284:ALA:HB3	17:AR:286:GLU:CD	2.39	0.42
17:AR:33:LEU:N	17:AR:33:LEU:CD2	2.82	0.42
1:AA:1573:G:N2	19:A7:41:U:C5'	2.39	0.42
1:AA:1461:C:H4'	19:A7:31:A:O5'	2.20	0.42
17:AR:133:VAL:CG1	17:AR:142:ALA:N	2.82	0.42
53:B5:271:C:C5	53:B5:272:G:H1'	2.55	0.42
23:B8:4:ILE:HG12	23:B8:6:GLU:H	1.84	0.42
17:AR:118:LYS:O	17:AR:119:ALA:HB3	2.20	0.42
17:AR:106:HIS:NE2	17:AR:132:LYS:HD2	2.35	0.42
32:BH:40:HIS:CD2	32:BH:41:ILE:H	2.37	0.42
48:BX:81:ALA:HA	53:B5:17:G:C4	2.54	0.42
17:AR:288:HIS:O	17:AR:306:THR:HG23	2.19	0.42
17:AR:147:HIS:CD2	17:AR:151:VAL:HG22	2.55	0.42
28:BD:208:VAL:HG11	28:BD:250:TRP:CZ3	2.55	0.42
29:BE:193:GLU:H	29:BE:213:ASP:H	1.68	0.42
53:B5:126:U:H2'	53:B5:127:G:C8	2.55	0.42
27:BC:32:PHE:CD1	53:B5:3003:G:H4'	2.53	0.42
17:AR:285:ALA:C	17:AR:305:TYR:HE2	2.22	0.42
17:AR:29:GLN:N	17:AR:30:PRO:CD	2.83	0.42
1:AA:115:G:H1	1:AA:302:U:H1'	1.85	0.42
17:AR:96:THR:CG2	17:AR:97:GLY:N	2.82	0.42
17:AR:164:ASP:C	17:AR:183:LEU:H	2.21	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1574:A:C6	19:A7:41:U:H4'	2.54	0.42
17:AR:139:GLN:HA	17:AR:139:GLN:OE1	2.20	0.42
17:AR:260:ILE:HG13	17:AR:284:ALA:HB2	1.98	0.42
17:AR:313:TRP:O	17:AR:314:GLN:HB2	2.20	0.42
17:AR:33:LEU:HD21	17:AR:45:TRP:CB	2.50	0.42
17:AR:135:THR:N	17:AR:141:LEU:CD1	2.83	0.42
17:AR:165:ASP:N	17:AR:184:ASN:N	2.68	0.42
17:AR:165:ASP:CB	17:AR:183:LEU:HB3	2.50	0.42
17:AR:203:THR:CG2	17:AR:245:PHE:CE1	3.01	0.42
1:AA:1520:U:C5'	1:AA:1521:G:P	3.05	0.41
1:AA:1657:A:H4'	42:BR:67:PRO:CG	2.47	0.41
17:AR:172:ALA:CB	17:AR:199:ILE:CD1	2.94	0.41
1:AA:865:A:H2'	1:AA:866:G:OP2	2.19	0.41
1:AA:703:G:H4'	40:BP:176:ALA:HB2	2.02	0.41
17:AR:245:PHE:HE2	17:AR:265:LEU:HD11	1.84	0.41
37:BM:178:ALA:O	37:BM:181:ALA:HB3	2.21	0.41
1:AA:1755:G:H2'	53:B5:2262:A:C2	2.54	0.41
1:AA:1520:U:C6	1:AA:1520:U:H3'	2.55	0.41
1:AA:1502:G:O6	1:AA:1547:C:C6	2.72	0.41
17:AR:135:THR:N	17:AR:141:LEU:HD11	2.34	0.41
17:AR:157:VAL:O	17:AR:157:VAL:HG13	2.20	0.41
17:AR:262:VAL:CG2	17:AR:272:ASP:HB3	2.49	0.41
17:AR:224:ASN:OD1	17:AR:227:ALA:HB3	2.20	0.41
42:BR:48:ARG:HA	53:B5:2338:C:H4'	2.01	0.41
17:AR:79:TYR:HD2	17:AR:92:TRP:O	2.03	0.41
1:AA:1120:C:O2'	53:B5:2190:U:C4'	2.68	0.41
1:AA:1775:G:C5'	53:B5:2193:U:OP1	2.68	0.41
17:AR:13:LEU:CD2	17:AR:13:LEU:N	2.83	0.41
17:AR:243:LEU:HD23	17:AR:254:ALA:HA	2.01	0.41
53:B5:2621:G:C8	53:B5:2622:C:C5	3.09	0.41
1:AA:501:U:C1'	1:AA:501:U:C6	2.93	0.41
14:AO:149:LEU:HD13	53:B5:847:A:P	2.46	0.41
17:AR:278:PHE:CZ	17:AR:281:TYR:O	2.73	0.41
17:AR:300:THR:HG22	17:AR:301:LEU:N	2.36	0.41
17:AR:131:ILE:HD12	17:AR:154:VAL:HG11	2.01	0.41
17:AR:233:THR:O	17:AR:234:LEU:HD23	2.21	0.41
1:AA:223:U:H3	1:AA:244:A:H61	1.68	0.41
12:AM:92:ILE:H	12:AM:92:ILE:HD13	1.84	0.41
53:B5:1384:U:H1'	53:B5:1385:C:C2	2.55	0.41
9:AJ:56:VAL:HG22	9:AJ:57:ARG:N	2.36	0.41
1:AA:628:G:H4'	53:B5:846:A:N3	2.36	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
17:AR:256:THR:O	17:AR:283:LYS:CG	2.66	0.41
17:AR:305:TYR:CE2	17:AR:311:ARG:NE	2.85	0.41
17:AR:315:VAL:CG1	17:AR:316:MET:N	2.83	0.41
17:AR:136:ILE:CG2	17:AR:137:LYS:N	2.83	0.41
17:AR:157:VAL:CG1	17:AR:167:VAL:HB	2.51	0.41
17:AR:267:PRO:O	17:AR:268:GLN:HB2	2.20	0.41
2:AB:23:HIS:HA	2:AB:48:ILE:HG22	2.03	0.41
1:AA:799:A:O3'	40:BP:160:ALA:O	2.39	0.41
1:AA:921:G:O2'	26:BB:141:PRO:HG3	2.20	0.41
17:AR:170:ILE:N	17:AR:170:ILE:CD1	2.82	0.41
8:AI:93:HIS:CE1	17:AR:59:ARG:CZ	3.03	0.41
8:AI:97:VAL:HG12	8:AI:98:ASP:N	2.36	0.41
6:AG:77:TYR:H	6:AG:191:ALA:HB2	1.85	0.41
19:A7:21:A:C6	19:A7:48:C:C6	3.09	0.41
4:AD:155:HIS:CG	4:AD:156:ILE:H	2.39	0.41
1:AA:1573:G:N3	19:A7:41:U:C5'	2.67	0.41
1:AA:921:G:O3'	26:BB:140:ASN:ND2	2.54	0.41
17:AR:276:PRO:N	17:AR:284:ALA:O	2.54	0.41
17:AR:72:THR:CG2	17:AR:73:LEU:N	2.82	0.41
17:AR:219:GLU:HG2	17:AR:220:ILE:N	2.36	0.41
43:BS:49:ILE:H	43:BS:49:ILE:HD12	1.86	0.41
37:BM:150:ALA:C	37:BM:154:ALA:HB3	2.41	0.41
1:AA:1120:C:HO2'	53:B5:2190:U:C4'	2.34	0.41
19:A7:10:2MG:HM23	19:A7:11:C:H1'	2.03	0.41
1:AA:856:A:C2	7:AH:33:VAL:CA	3.04	0.40
1:AA:1430:G:H1'	13:AN:38:ILE:HD11	2.03	0.40
28:BD:55:LYS:HA	28:BD:58:HIS:CD2	2.56	0.40
36:BL:80:THR:HG23	53:B5:2607:G:H22	1.86	0.40
17:AR:286:GLU:OE2	17:AR:305:TYR:CE1	2.74	0.40
17:AR:303:ALA:O	17:AR:305:TYR:CE1	2.74	0.40
17:AR:13:LEU:HD23	17:AR:310:ILE:HG22	2.03	0.40
17:AR:61:PHE:HE2	17:AR:97:GLY:HA2	1.73	0.40
46:BV:74:LEU:HD13	46:BV:74:LEU:H	1.86	0.40
17:AR:37:SER:HB3	17:AR:39:ASP:OD1	2.21	0.40
25:BA:105:LYS:HE2	53:B5:2477:G:C5	2.57	0.40
17:AR:285:ALA:HB2	17:AR:313:TRP:CH2	2.57	0.40
17:AR:172:ALA:HB1	17:AR:199:ILE:HG23	2.02	0.40
17:AR:183:LEU:HA	17:AR:183:LEU:HD13	1.98	0.40
1:AA:701:U:C2	40:BP:171:ALA:C	2.91	0.40
1:AA:1744:A:C5'	53:B5:2291:A:H1'	2.47	0.40
17:AR:286:GLU:CD	17:AR:305:TYR:CE2	2.95	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
17:AR:285:ALA:HB2	17:AR:313:TRP:HH2	1.86	0.40
17:AR:68:VAL:HG21	17:AR:82:SER:HB2	2.04	0.40
17:AR:42:LEU:CD2	17:AR:71:CYS:HB2	2.52	0.40
17:AR:34:LEU:CD1	17:AR:73:LEU:HD21	2.48	0.40
17:AR:79:TYR:CD2	17:AR:93:ASP:HA	2.55	0.40
17:AR:199:ILE:CD1	17:AR:202:LEU:CD2	2.97	0.40
17:AR:133:VAL:HG12	17:AR:142:ALA:H	1.87	0.40
17:AR:234:LEU:HD13	17:AR:263:PHE:CD2	2.57	0.40
11:AL:62:LYS:HG2	11:AL:63:GLN:H	1.86	0.40
9:AJ:30:LYS:HG3	9:AJ:31:VAL:H	1.85	0.40
37:BM:143:THR:HG23	37:BM:144:SER:H	1.86	0.40
1:AA:858:G:H2'	7:AH:28:ARG:N	2.37	0.40
17:AR:106:HIS:CD2	17:AR:132:LYS:HD2	2.56	0.40
17:AR:262:VAL:CG2	17:AR:272:ASP:H	2.30	0.40
23:B8:24:SER:HA	23:B8:115:ALA:H	1.86	0.40
23:B8:84:VAL:HG23	23:B8:110:ARG:HE	1.87	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 PDB entries followed by that with respect to all EM entries.

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	191/193 (99%)	173 (91%)	16 (8%)	2 (1%)	19	65
3	AC	186/188 (99%)	148 (80%)	28 (15%)	10 (5%)	2	29
4	AD	120/158 (76%)	97 (81%)	16 (13%)	7 (6%)	2	27
5	AE	160/162 (99%)	141 (88%)	13 (8%)	6 (4%)	4	37
6	AG	184/186 (99%)	156 (85%)	20 (11%)	8 (4%)	3	34
7	AH	121/125 (97%)	105 (87%)	11 (9%)	5 (4%)	3	35
8	AI	136/138 (99%)	106 (78%)	21 (15%)	9 (7%)	1	24

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
9	AJ	94/96 (98%)	77 (82%)	14 (15%)	3 (3%)	5	41
10	AK	123/125 (98%)	99 (80%)	20 (16%)	4 (3%)	5	40
11	AL	116/118 (98%)	89 (77%)	25 (22%)	2 (2%)	11	55
12	AM	128/130 (98%)	103 (80%)	20 (16%)	5 (4%)	4	36
13	AN	48/50 (96%)	36 (75%)	8 (17%)	4 (8%)	1	18
14	AO	82/84 (98%)	68 (83%)	11 (13%)	3 (4%)	4	38
15	AQ	78/80 (98%)	66 (85%)	11 (14%)	1 (1%)	15	60
16	AS	69/71 (97%)	54 (78%)	12 (17%)	3 (4%)	3	34
17	AR	311/313 (99%)	282 (91%)	28 (9%)	1 (0%)	46	83
18	AT	137/141 (97%)	110 (80%)	21 (15%)	6 (4%)	3	33
20	B0	107/109 (98%)	75 (70%)	20 (19%)	12 (11%)	0	11
21	B1	46/48 (96%)	36 (78%)	9 (20%)	1 (2%)	8	49
22	B2	96/98 (98%)	87 (91%)	7 (7%)	2 (2%)	9	50
23	B8	116/118 (98%)	85 (73%)	25 (22%)	6 (5%)	2	30
24	B9	70/72 (97%)	48 (69%)	16 (23%)	6 (9%)	1	17
25	BA	211/213 (99%)	172 (82%)	30 (14%)	9 (4%)	3	34
26	BB	241/243 (99%)	174 (72%)	53 (22%)	14 (6%)	2	27
27	BC	360/362 (99%)	279 (78%)	68 (19%)	13 (4%)	4	38
28	BD	255/257 (99%)	208 (82%)	41 (16%)	6 (2%)	7	47
29	BE	235/237 (99%)	175 (74%)	47 (20%)	13 (6%)	2	29
30	BF	211/213 (99%)	150 (71%)	55 (26%)	6 (3%)	6	44
31	BG	111/113 (98%)	88 (79%)	20 (18%)	3 (3%)	6	45
32	BH	177/179 (99%)	146 (82%)	25 (14%)	6 (3%)	5	40
33	BI	163/165 (99%)	123 (76%)	31 (19%)	9 (6%)	2	29
34	BJ	149/151 (99%)	112 (75%)	33 (22%)	4 (3%)	6	45
35	BK	136/138 (99%)	107 (79%)	22 (16%)	7 (5%)	2	30
36	BL	190/192 (99%)	156 (82%)	29 (15%)	5 (3%)	7	45
37	BM	176/178 (99%)	134 (76%)	34 (19%)	8 (4%)	3	33
38	BN	148/150 (99%)	122 (82%)	23 (16%)	3 (2%)	9	51
39	BO	119/121 (98%)	92 (77%)	23 (19%)	4 (3%)	5	40
40	BP	174/176 (99%)	155 (89%)	15 (9%)	4 (2%)	8	48

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
41	BQ	114/116 (98%)	94 (82%)	16 (14%)	4 (4%)	4	39
42	BR	129/131 (98%)	96 (74%)	28 (22%)	5 (4%)	4	36
43	BS	43/45 (96%)	33 (77%)	10 (23%)	0	100	100
44	BT	78/80 (98%)	59 (76%)	16 (20%)	3 (4%)	4	37
45	BU	114/116 (98%)	95 (83%)	15 (13%)	4 (4%)	4	39
46	BV	140/142 (99%)	98 (70%)	31 (22%)	11 (8%)	1	19
47	BW	77/79 (98%)	58 (75%)	16 (21%)	3 (4%)	4	36
48	BX	84/86 (98%)	63 (75%)	14 (17%)	7 (8%)	1	18
49	BY	50/52 (96%)	39 (78%)	10 (20%)	1 (2%)	9	51
50	BZ	90/92 (98%)	68 (76%)	18 (20%)	4 (4%)	3	33
All	All	6694/6830 (98%)	5337 (80%)	1095 (16%)	262 (4%)	7	36

All (262) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	AC	81	PRO
3	AC	179	GLN
4	AD	96	VAL
5	AE	86	VAL
5	AE	226	THR
5	AE	227	PRO
6	AG	89	ILE
6	AG	164	PRO
7	AH	35	ILE
8	AI	97	VAL
8	AI	134	ALA
8	AI	136	SER
9	AJ	72	ASN
11	AL	53	VAL
12	AM	51	ASP
13	AN	16	LYS
14	AO	108	ASP
16	AS	104	GLN
18	AT	51	GLU
20	B0	38	ILE
20	B0	50	ILE
20	B0	64	LYS
24	B9	14	TYR

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Mol	Chain	Res	Type
24	B9	34	HIS
25	BA	60	ARG
26	BB	140	ASN
26	BB	161	ASP
26	BB	197	PRO
27	BC	168	LYS
27	BC	279	ASN
29	BE	188	GLU
29	BE	212	ALA
29	BE	216	GLU
30	BF	83	LEU
30	BF	123	THR
33	BI	18	PRO
33	BI	104	SER
35	BK	78	SER
35	BK	92	ARG
36	BL	56	LYS
37	BM	149	ALA
37	BM	170	ALA
46	BV	41	ARG
48	BX	68	ALA
48	BX	70	ALA
48	BX	80	ALA
50	BZ	46	LYS
50	BZ	57	VAL
2	AB	43	ASP
2	AB	68	PRO
3	AC	163	PRO
3	AC	164	VAL
4	AD	18	PRO
4	AD	148	VAL
5	AE	205	ARG
7	AH	58	SER
7	AH	98	GLN
8	AI	15	SER
11	AL	144	ARG
12	AM	137	HIS
14	AO	128	TYR
18	AT	37	VAL
20	B0	77	ALA
20	B0	116	GLY
23	B8	35	SER

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Mol	Chain	Res	Type
25	BA	36	VAL
25	BA	105	LYS
25	BA	162	VAL
26	BB	156	LYS
27	BC	128	LYS
27	BC	129	ALA
28	BD	180	LYS
29	BE	167	SER
29	BE	183	TRP
29	BE	224	LYS
31	BG	142	LEU
33	BI	142	ASP
35	BK	10	VAL
36	BL	40	ALA
37	BM	60	LYS
37	BM	156	ALA
40	BP	88	ARG
40	BP	143	ALA
42	BR	37	ILE
46	BV	52	PHE
46	BV	101	ILE
47	BW	61	LYS
48	BX	86	ALA
3	AC	83	THR
3	AC	112	GLY
3	AC	145	ALA
3	AC	148	LYS
4	AD	9	SER
4	AD	87	SER
4	AD	88	GLU
6	AG	66	GLN
6	AG	75	GLY
6	AG	85	ALA
6	AG	163	SER
7	AH	28	ARG
7	AH	84	GLY
8	AI	42	GLU
8	AI	139	GLN
17	AR	283	LYS
18	AT	53	TRP
18	AT	86	ARG
20	B0	48	GLY

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Mol	Chain	Res	Type
20	B0	62	LYS
20	B0	65	PHE
21	B1	40	LYS
23	B8	89	THR
24	B9	59	CYS
24	B9	66	GLY
26	BB	70	ARG
27	BC	38	SER
27	BC	46	PHE
27	BC	365	PHE
29	BE	36	LEU
29	BE	194	LEU
29	BE	219	PHE
30	BF	120	THR
31	BG	191	ASN
32	BH	22	SER
33	BI	105	CYS
33	BI	144	ASN
34	BJ	147	THR
35	BK	20	GLY
35	BK	88	PRO
37	BM	169	ALA
38	BN	122	ALA
39	BO	118	GLY
39	BO	122	ILE
40	BP	73	GLY
40	BP	133	LYS
41	BQ	10	ARG
44	BT	62	VAL
46	BV	14	HIS
47	BW	78	LYS
48	BX	48	ARG
50	BZ	35	LEU
3	AC	9	ARG
4	AD	19	TYR
8	AI	125	GLU
10	AK	97	GLY
12	AM	90	ASN
13	AN	39	CYS
14	AO	147	SER
20	B0	94	ALA
22	B2	82	GLY

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Mol	Chain	Res	Type
23	B8	116	PRO
24	B9	61	LYS
25	BA	58	CYS
25	BA	196	LYS
25	BA	208	SER
26	BB	73	GLU
26	BB	129	ALA
26	BB	153	GLY
26	BB	179	LEU
26	BB	221	LYS
26	BB	230	VAL
27	BC	102	LEU
27	BC	234	GLY
27	BC	272	TYR
28	BD	186	LYS
28	BD	213	ASN
28	BD	242	ALA
29	BE	227	LEU
30	BF	30	ALA
30	BF	177	GLY
30	BF	226	GLY
31	BG	126	SER
32	BH	129	ARG
33	BI	25	ALA
34	BJ	4	LYS
35	BK	26	ALA
37	BM	37	ARG
39	BO	74	GLU
41	BQ	18	ASP
41	BQ	69	LYS
42	BR	72	LYS
42	BR	96	GLU
44	BT	104	GLU
45	BU	59	VAL
46	BV	63	GLN
46	BV	114	LYS
47	BW	83	GLU
48	BX	38	ARG
48	BX	82	ALA
3	AC	152	PHE
6	AG	86	GLN
8	AI	32	ASN

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Mol	Chain	Res	Type
9	AJ	97	VAL
10	AK	47	LYS
10	AK	55	SER
13	AN	37	ASN
16	AS	51	SER
18	AT	89	ARG
23	B8	32	ASN
23	B8	46	ARG
26	BB	33	ASP
26	BB	180	LEU
28	BD	15	ALA
32	BH	50	ASN
32	BH	118	LEU
32	BH	122	LYS
33	BI	101	LYS
33	BI	111	LEU
41	BQ	17	ARG
42	BR	21	ALA
45	BU	46	LYS
46	BV	32	GLY
46	BV	102	ASP
46	BV	141	GLY
49	BY	36	SER
8	AI	126	PRO
18	AT	90	PRO
20	B0	97	ALA
25	BA	125	GLY
27	BC	29	VAL
29	BE	87	GLY
29	BE	201	GLY
29	BE	223	PHE
32	BH	106	LYS
34	BJ	12	LEU
35	BK	86	LYS
36	BL	173	GLY
36	BL	191	TRP
37	BM	58	LEU
38	BN	65	SER
42	BR	44	SER
45	BU	113	LYS
5	AE	145	GLY
12	AM	75	ASN

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Mol	Chain	Res	Type
15	AQ	111	VAL
16	AS	99	GLY
20	B0	96	ILE
24	B9	43	GLY
25	BA	37	GLY
34	BJ	58	GLY
39	BO	96	PHE
44	BT	86	VAL
46	BV	143	VAL
50	BZ	56	PRO
6	AG	162	VAL
9	AJ	91	ILE
22	B2	81	VAL
27	BC	260	VAL
37	BM	30	GLY
5	AE	109	GLY
12	AM	95	GLY
28	BD	139	GLY
38	BN	19	GLY
45	BU	44	GLY
46	BV	15	VAL
10	AK	35	GLY
13	AN	29	GLY
23	B8	70	LEU
26	BB	121	GLY
33	BI	114	GLY
36	BL	52	GLY
20	B0	79	VAL
27	BC	307	PRO

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all EM entries.

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	161/161 (100%)	160 (99%)	1 (1%)	90	95
3	AC	152/152 (100%)	149 (98%)	3 (2%)	63	85

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
4	AD	113/142 (80%)	103 (91%)	10 (9%)	12	45
5	AE	127/127 (100%)	122 (96%)	5 (4%)	39	72
6	AG	155/155 (100%)	150 (97%)	5 (3%)	46	76
7	AH	106/106 (100%)	100 (94%)	6 (6%)	25	62
8	AI	115/115 (100%)	107 (93%)	8 (7%)	19	56
9	AJ	90/90 (100%)	86 (96%)	4 (4%)	35	69
10	AK	95/95 (100%)	88 (93%)	7 (7%)	17	54
11	AL	98/98 (100%)	94 (96%)	4 (4%)	37	71
12	AM	115/115 (100%)	107 (93%)	8 (7%)	19	56
13	AN	45/45 (100%)	40 (89%)	5 (11%)	8	34
14	AO	74/74 (100%)	70 (95%)	4 (5%)	27	64
15	AQ	71/71 (100%)	68 (96%)	3 (4%)	36	70
16	AS	57/57 (100%)	56 (98%)	1 (2%)	66	87
17	AR	257/257 (100%)	246 (96%)	11 (4%)	35	70
18	AT	114/114 (100%)	102 (90%)	12 (10%)	8	36
20	B0	94/94 (100%)	89 (95%)	5 (5%)	28	64
21	B1	44/44 (100%)	41 (93%)	3 (7%)	20	57
22	B2	82/82 (100%)	80 (98%)	2 (2%)	57	82
23	B8	100/100 (100%)	87 (87%)	13 (13%)	5	28
24	B9	55/55 (100%)	52 (94%)	3 (6%)	27	63
25	BA	194/194 (100%)	183 (94%)	11 (6%)	25	62
26	BB	186/186 (100%)	176 (95%)	10 (5%)	27	64
27	BC	302/302 (100%)	282 (93%)	20 (7%)	21	57
28	BD	199/199 (100%)	188 (94%)	11 (6%)	27	63
29	BE	199/199 (100%)	187 (94%)	12 (6%)	24	60
30	BF	143/143 (100%)	135 (94%)	8 (6%)	26	62
31	BG	89/89 (100%)	82 (92%)	7 (8%)	15	51
32	BH	160/160 (100%)	151 (94%)	9 (6%)	26	62
33	BI	141/141 (100%)	134 (95%)	7 (5%)	30	66
34	BJ	129/129 (100%)	120 (93%)	9 (7%)	19	56
35	BK	112/112 (100%)	107 (96%)	5 (4%)	34	69

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
36	BL	164/164 (100%)	157 (96%)	7 (4%)	35	70
37	BM	119/119 (100%)	108 (91%)	11 (9%)	11	43
38	BN	122/122 (100%)	118 (97%)	4 (3%)	45	76
39	BO	99/99 (100%)	96 (97%)	3 (3%)	48	77
40	BP	118/118 (100%)	113 (96%)	5 (4%)	36	70
41	BQ	87/87 (100%)	81 (93%)	6 (7%)	19	56
42	BR	102/102 (100%)	96 (94%)	6 (6%)	24	61
43	BS	38/38 (100%)	32 (84%)	6 (16%)	3	21
44	BT	71/71 (100%)	64 (90%)	7 (10%)	10	39
45	BU	100/100 (100%)	98 (98%)	2 (2%)	63	85
46	BV	113/113 (100%)	103 (91%)	10 (9%)	12	45
47	BW	69/69 (100%)	65 (94%)	4 (6%)	25	61
48	BX	55/55 (100%)	53 (96%)	2 (4%)	42	74
49	BY	42/42 (100%)	39 (93%)	3 (7%)	18	55
50	BZ	81/81 (100%)	76 (94%)	5 (6%)	23	60
All	All	5554/5583 (100%)	5241 (94%)	313 (6%)	31	62

All (313) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	AB	179	ARG
3	AC	7	LYS
3	AC	132	LYS
3	AC	143	ARG
4	AD	8	TYR
4	AD	111	THR
4	AD	112	GLN
4	AD	120	LYS
4	AD	127	VAL
4	AD	138	LYS
4	AD	139	GLN
4	AD	146	PHE
4	AD	153	GLU
4	AD	154	LYS
5	AE	82	ASN
5	AE	111	VAL
5	AE	139	ILE

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Mol	Chain	Res	Type
5	AE	178	ILE
5	AE	204	THR
6	AG	106	LYS
6	AG	156	ARG
6	AG	180	ARG
6	AG	219	ARG
6	AG	224	ASN
7	AH	9	ASP
7	AH	27	ILE
7	AH	32	LYS
7	AH	35	ILE
7	AH	37	PHE
7	AH	50	PHE
8	AI	8	GLN
8	AI	43	ILE
8	AI	45	ARG
8	AI	89	LEU
8	AI	96	TYR
8	AI	98	ASP
8	AI	127	LYS
8	AI	142	TYR
9	AJ	21	LYS
9	AJ	43	LYS
9	AJ	74	GLU
9	AJ	91	ILE
10	AK	26	THR
10	AK	33	LEU
10	AK	39	ILE
10	AK	47	LYS
10	AK	66	ASP
10	AK	115	ILE
10	AK	128	LYS
11	AL	117	ILE
11	AL	136	TRP
11	AL	140	LYS
11	AL	141	GLU
12	AM	18	LEU
12	AM	54	LEU
12	AM	57	ARG
12	AM	66	LEU
12	AM	92	ILE
12	AM	116	LEU

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Mol	Chain	Res	Type
12	AM	125	ILE
12	AM	132	ARG
13	AN	14	PHE
13	AN	22	ARG
13	AN	31	VAL
13	AN	39	CYS
13	AN	44	ARG
14	AO	78	ASN
14	AO	114	ARG
14	AO	125	LEU
14	AO	134	VAL
15	AQ	67	ARG
15	AQ	80	MET
15	AQ	131	ILE
16	AS	97	TYR
17	AR	13	LEU
17	AR	32	LEU
17	AR	33	LEU
17	AR	52	GLN
17	AR	118	LYS
17	AR	193	ILE
17	AR	199	ILE
17	AR	228	LYS
17	AR	290	VAL
17	AR	292	LEU
17	AR	308	ASN
18	AT	6	VAL
18	AT	7	ARG
18	AT	24	ARG
18	AT	29	GLU
18	AT	77	ASN
18	AT	86	ARG
18	AT	89	ARG
18	AT	92	LYS
18	AT	94	ILE
18	AT	103	LYS
18	AT	123	ARG
18	AT	126	GLU
20	B0	21	HIS
20	B0	38	ILE
20	B0	39	ASP
20	B0	65	PHE

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Mol	Chain	Res	Type
20	B0	111	ARG
21	B1	9	ILE
21	B1	20	ASN
21	B1	49	MET
22	B2	28	LYS
22	B2	104	LEU
23	B8	14	LYS
23	B8	33	VAL
23	B8	36	GLN
23	B8	38	MET
23	B8	39	HIS
23	B8	43	LYS
23	B8	48	ARG
23	B8	51	VAL
23	B8	64	ARG
23	B8	75	LYS
23	B8	97	LYS
23	B8	98	ASN
23	B8	116	PRO
24	B9	22	LEU
24	B9	41	PHE
24	B9	45	LYS
25	BA	4	ILE
25	BA	12	HIS
25	BA	29	LEU
25	BA	31	THR
25	BA	45	ARG
25	BA	61	PRO
25	BA	76	ARG
25	BA	162	VAL
25	BA	199	GLN
25	BA	200	ASN
25	BA	214	PHE
26	BB	5	ILE
26	BB	42	ARG
26	BB	61	VAL
26	BB	64	ARG
26	BB	65	ASP
26	BB	75	ILE
26	BB	145	LYS
26	BB	155	LYS
26	BB	181	LYS

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Mol	Chain	Res	Type
26	BB	219	ILE
27	BC	20	LYS
27	BC	32	PHE
27	BC	37	ARG
27	BC	59	ASP
27	BC	121	ASN
27	BC	126	LYS
27	BC	127	LYS
27	BC	150	ARG
27	BC	153	LYS
27	BC	168	LYS
27	BC	215	ILE
27	BC	221	THR
27	BC	237	LYS
27	BC	240	ARG
27	BC	244	ARG
27	BC	317	ILE
27	BC	327	CYS
27	BC	332	ARG
27	BC	360	ASP
27	BC	364	LYS
28	BD	33	ASP
28	BD	50	TYR
28	BD	69	ARG
28	BD	113	VAL
28	BD	119	ARG
28	BD	126	ILE
28	BD	148	ILE
28	BD	170	LYS
28	BD	191	LYS
28	BD	194	TYR
28	BD	215	ILE
29	BE	23	ARG
29	BE	48	LYS
29	BE	50	ARG
29	BE	105	ILE
29	BE	131	LEU
29	BE	151	GLN
29	BE	173	VAL
29	BE	196	ARG
29	BE	213	ASP
29	BE	217	GLU

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Mol	Chain	Res	Type
29	BE	219	PHE
29	BE	247	ILE
30	BF	83	LEU
30	BF	84	VAL
30	BF	111	ILE
30	BF	118	LYS
30	BF	125	GLU
30	BF	129	LEU
30	BF	185	ILE
30	BF	244	ASN
31	BG	107	GLU
31	BG	134	TYR
31	BG	136	LEU
31	BG	146	LYS
31	BG	163	VAL
31	BG	172	LYS
31	BG	200	LEU
32	BH	11	GLU
32	BH	79	ILE
32	BH	116	ASN
32	BH	149	ASN
32	BH	167	VAL
32	BH	170	LYS
32	BH	172	ILE
32	BH	177	ASP
32	BH	183	HIS
33	BI	69	ARG
33	BI	82	ARG
33	BI	103	LEU
33	BI	115	MET
33	BI	121	LYS
33	BI	134	ILE
33	BI	159	PHE
34	BJ	10	ARG
34	BJ	44	THR
34	BJ	47	GLN
34	BJ	52	TYR
34	BJ	78	GLU
34	BJ	94	ARG
34	BJ	123	PHE
34	BJ	131	MET
34	BJ	137	ARG

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Mol	Chain	Res	Type
35	BK	18	VAL
35	BK	60	VAL
35	BK	87	GLU
35	BK	93	LYS
35	BK	96	LYS
36	BL	13	LYS
36	BL	56	LYS
36	BL	67	ARG
36	BL	77	LYS
36	BL	86	ASN
36	BL	122	ASN
36	BL	191	TRP
37	BM	14	HIS
37	BM	39	GLU
37	BM	42	ASN
37	BM	50	ASN
37	BM	58	LEU
37	BM	84	LEU
37	BM	96	LYS
37	BM	113	ASP
37	BM	114	LYS
37	BM	134	LYS
37	BM	140	LYS
38	BN	69	ARG
38	BN	118	GLN
38	BN	124	LYS
38	BN	153	LYS
39	BO	48	VAL
39	BO	72	LYS
39	BO	98	LYS
40	BP	20	ARG
40	BP	63	THR
40	BP	96	ILE
40	BP	111	ASP
40	BP	125	LYS
41	BQ	26	HIS
41	BQ	27	LEU
41	BQ	35	LYS
41	BQ	78	LYS
41	BQ	84	TYR
41	BQ	91	LEU
42	BR	11	PHE

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Mol	Chain	Res	Type
42	BR	36	ILE
42	BR	59	MET
42	BR	66	LYS
42	BR	88	ARG
42	BR	94	TYR
43	BS	14	TYR
43	BS	21	PHE
43	BS	28	ILE
43	BS	30	ARG
43	BS	43	ARG
43	BS	48	ARG
44	BT	60	TYR
44	BT	61	LYS
44	BT	67	ILE
44	BT	78	ASP
44	BT	89	LYS
44	BT	93	TYR
44	BT	95	ILE
45	BU	59	VAL
45	BU	97	ILE
46	BV	47	TYR
46	BV	62	LYS
46	BV	69	LYS
46	BV	74	LEU
46	BV	91	LYS
46	BV	97	THR
46	BV	127	ARG
46	BV	131	LYS
46	BV	137	ILE
46	BV	147	ILE
47	BW	11	GLU
47	BW	61	LYS
47	BW	70	ARG
47	BW	83	GLU
48	BX	43	LYS
48	BX	45	LYS
49	BY	12	HIS
49	BY	17	THR
49	BY	43	LYS
50	BZ	6	LYS
50	BZ	9	LYS
50	BZ	38	GLN

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Mol	Chain	Res	Type
50	BZ	40	LYS
50	BZ	83	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (47) such sidechains are listed below:

Mol	Chain	Res	Type
2	AB	23	HIS
3	AC	159	HIS
4	AD	124	HIS
5	AE	233	GLN
6	AG	66	GLN
6	AG	72	HIS
8	AI	74	HIS
8	AI	93	HIS
12	AM	78	HIS
12	AM	122	HIS
15	AQ	110	HIS
16	AS	79	HIS
16	AS	98	ASN
17	AR	31	ASN
17	AR	52	GLN
17	AR	101	GLN
17	AR	185	GLN
17	AR	198	ASN
20	B0	71	HIS
25	BA	12	HIS
25	BA	44	GLN
27	BC	173	GLN
27	BC	177	HIS
27	BC	313	HIS
28	BD	36	HIS
28	BD	58	HIS
30	BF	112	ASN
30	BF	194	HIS
30	BF	231	ASN
31	BG	191	ASN
32	BH	40	HIS
32	BH	116	ASN
32	BH	157	ASN
33	BI	92	HIS
33	BI	95	HIS
34	BJ	47	GLN

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Mol	Chain	Res	Type
36	BL	123	GLN
36	BL	138	GLN
36	BL	156	HIS
37	BM	14	HIS
39	BO	135	GLN
41	BQ	22	HIS
46	BV	14	HIS
46	BV	39	HIS
46	BV	119	ASN
49	BY	16	HIS
50	BZ	23	HIS

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	AA	1757/1761 (99%)	380 (21%)	79 (4%)
19	A7	75/76 (98%)	8 (10%)	3 (4%)
51	B3	112/113 (99%)	40 (35%)	4 (3%)
52	B4	156/157 (99%)	55 (35%)	10 (6%)
53	B5	3169/3170 (99%)	942 (29%)	139 (4%)
All	All	5269/5277 (99%)	1425 (27%)	235 (4%)

All (1425) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	AA	2	A
1	AA	3	U
1	AA	4	C
1	AA	9	U
1	AA	27	U
1	AA	34	G
1	AA	42	G
1	AA	44	U
1	AA	46	A
1	AA	47	A
1	AA	48	G
1	AA	50	C
1	AA	51	A
1	AA	57	G
1	AA	60	U
1	AA	61	A

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Mol	Chain	Res	Type
1	AA	62	A
1	AA	65	A
1	AA	67	A
1	AA	68	A
1	AA	69	G
1	AA	81	G
1	AA	82	U
1	AA	83	G
1	AA	86	A
1	AA	90	C
1	AA	92	A
1	AA	100	A
1	AA	105	A
1	AA	106	U
1	AA	109	G
1	AA	110	U
1	AA	115	G
1	AA	116	U
1	AA	147	A
1	AA	161	U
1	AA	174	U
1	AA	176	C
1	AA	180	A
1	AA	188	A
1	AA	193	U
1	AA	200	A
1	AA	204	G
1	AA	208	U
1	AA	213	A
1	AA	230	C
1	AA	232	U
1	AA	233	C
1	AA	241	U
1	AA	243	G
1	AA	256	A
1	AA	257	A
1	AA	263	C
1	AA	264	G
1	AA	265	A
1	AA	266	A
1	AA	267	U
1	AA	279	G

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Mol	Chain	Res	Type
1	AA	280	U
1	AA	281	G
1	AA	282	C
1	AA	283	U
1	AA	284	G
1	AA	291	G
1	AA	295	A
1	AA	308	C
1	AA	309	C
1	AA	312	A
1	AA	313	U
1	AA	314	C
1	AA	316	A
1	AA	319	U
1	AA	322	G
1	AA	328	A
1	AA	329	G
1	AA	337	G
1	AA	351	C
1	AA	361	C
1	AA	365	G
1	AA	366	A
1	AA	373	G
1	AA	400	A
1	AA	401	A
1	AA	404	G
1	AA	419	G
1	AA	423	G
1	AA	424	C
1	AA	426	G
1	AA	429	G
1	AA	435	C
1	AA	439	U
1	AA	445	A
1	AA	451	A
1	AA	452	A
1	AA	462	G
1	AA	468	A
1	AA	469	C
1	AA	475	A
1	AA	478	A
1	AA	479	C

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Mol	Chain	Res	Type
1	AA	481	A
1	AA	493	U
1	AA	494	U
1	AA	495	C
1	AA	496	G
1	AA	501	U
1	AA	503	G
1	AA	512	A
1	AA	513	U
1	AA	530	C
1	AA	542	A
1	AA	543	C
1	AA	551	G
1	AA	558	U
1	AA	565	C
1	AA	568	G
1	AA	571	G
1	AA	574	G
1	AA	578	U
1	AA	579	A
1	AA	594	A
1	AA	609	U
1	AA	611	U
1	AA	613	G
1	AA	619	A
1	AA	620	A
1	AA	623	A
1	AA	624	G
1	AA	631	G
1	AA	638	U
1	AA	639	U
1	AA	648	G
1	AA	653	C
1	AA	666	U
1	AA	687	G
1	AA	693	U
1	AA	694	U
1	AA	695	U
1	AA	696	C
1	AA	698	U
1	AA	703	G
1	AA	729	G

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Mol	Chain	Res	Type
1	AA	730	G
1	AA	746	A
1	AA	749	U
1	AA	760	A
1	AA	765	G
1	AA	771	A
1	AA	774	A
1	AA	775	G
1	AA	776	G
1	AA	778	G
1	AA	779	U
1	AA	780	A
1	AA	781	U
1	AA	792	U
1	AA	797	G
1	AA	800	U
1	AA	818	C
1	AA	821	U
1	AA	822	U
1	AA	824	G
1	AA	844	A
1	AA	845	G
1	AA	853	G
1	AA	855	A
1	AA	856	A
1	AA	857	U
1	AA	858	G
1	AA	860	U
1	AA	864	U
1	AA	865	A
1	AA	866	G
1	AA	867	G
1	AA	868	G
1	AA	869	A
1	AA	877	G
1	AA	896	U
1	AA	898	G
1	AA	899	A
1	AA	900	G
1	AA	905	A
1	AA	923	A
1	AA	931	U

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Mol	Chain	Res	Type
1	AA	934	U
1	AA	957	U
1	AA	958	U
1	AA	961	C
1	AA	962	A
1	AA	963	U
1	AA	965	A
1	AA	971	G
1	AA	987	A
1	AA	989	C
1	AA	990	G
1	AA	993	G
1	AA	1003	U
1	AA	1004	A
1	AA	1022	A
1	AA	1025	A
1	AA	1027	C
1	AA	1029	A
1	AA	1031	G
1	AA	1036	C
1	AA	1038	A
1	AA	1039	G
1	AA	1040	A
1	AA	1041	U
1	AA	1042	C
1	AA	1049	U
1	AA	1050	G
1	AA	1056	U
1	AA	1058	A
1	AA	1059	A
1	AA	1060	U
1	AA	1074	C
1	AA	1090	A
1	AA	1092	U
1	AA	1098	G
1	AA	1099	G
1	AA	1110	A
1	AA	1127	G
1	AA	1134	A
1	AA	1135	A
1	AA	1148	A
1	AA	1152	G

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Mol	Chain	Res	Type
1	AA	1155	C
1	AA	1156	C
1	AA	1160	A
1	AA	1171	C
1	AA	1178	U
1	AA	1179	A
1	AA	1181	U
1	AA	1182	U
1	AA	1187	U
1	AA	1188	C
1	AA	1189	A
1	AA	1190	A
1	AA	1192	A
1	AA	1195	G
1	AA	1196	G
1	AA	1197	G
1	AA	1198	A
1	AA	1199	A
1	AA	1201	C
1	AA	1202	U
1	AA	1213	A
1	AA	1214	G
1	AA	1215	A
1	AA	1216	C
1	AA	1217	A
1	AA	1225	G
1	AA	1241	G
1	AA	1242	C
1	AA	1243	U
1	AA	1246	U
1	AA	1247	U
1	AA	1248	C
1	AA	1249	U
1	AA	1261	G
1	AA	1262	U
1	AA	1263	G
1	AA	1265	U
1	AA	1266	G
1	AA	1267	G
1	AA	1270	C
1	AA	1271	A
1	AA	1272	U

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Mol	Chain	Res	Type
1	AA	1274	G
1	AA	1281	C
1	AA	1282	U
1	AA	1318	A
1	AA	1319	A
1	AA	1325	G
1	AA	1329	C
1	AA	1341	A
1	AA	1342	A
1	AA	1350	U
1	AA	1351	G
1	AA	1355	G
1	AA	1366	U
1	AA	1378	U
1	AA	1379	A
1	AA	1386	C
1	AA	1387	U
1	AA	1396	C
1	AA	1402	G
1	AA	1409	G
1	AA	1410	U
1	AA	1411	U
1	AA	1416	G
1	AA	1424	A
1	AA	1425	G
1	AA	1430	G
1	AA	1431	U
1	AA	1438	C
1	AA	1443	A
1	AA	1444	A
1	AA	1445	C
1	AA	1446	G
1	AA	1447	U
1	AA	1452	G
1	AA	1455	C
1	AA	1456	G
1	AA	1457	C
1	AA	1458	A
1	AA	1459	C
1	AA	1471	U
1	AA	1472	G
1	AA	1479	C

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Mol	Chain	Res	Type
1	AA	1489	U
1	AA	1490	A
1	AA	1491	A
1	AA	1496	G
1	AA	1504	G
1	AA	1506	U
1	AA	1507	C
1	AA	1513	A
1	AA	1515	U
1	AA	1517	U
1	AA	1518	U
1	AA	1519	G
1	AA	1520	U
1	AA	1521	G
1	AA	1522	A
1	AA	1533	U
1	AA	1536	U
1	AA	1540	G
1	AA	1546	G
1	AA	1547	C
1	AA	1548	A
1	AA	1549	U
1	AA	1550	U
1	AA	1552	U
1	AA	1553	A
1	AA	1555	U
1	AA	1556	U
1	AA	1561	C
1	AA	1564	U
1	AA	1571	A
1	AA	1580	U
1	AA	1581	A
1	AA	1582	G
1	AA	1583	U
1	AA	1584	A
1	AA	1585	A
1	AA	1588	G
1	AA	1594	C
1	AA	1599	G
1	AA	1607	U
1	AA	1609	A
1	AA	1610	U

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Mol	Chain	Res	Type
1	AA	1612	A
1	AA	1613	C
1	AA	1614	G
1	AA	1615	U
1	AA	1631	A
1	AA	1633	A
1	AA	1634	C
1	AA	1635	C
1	AA	1636	G
1	AA	1637	C
1	AA	1638	C
1	AA	1640	G
1	AA	1655	U
1	AA	1656	G
1	AA	1676	A
1	AA	1753	A
1	AA	1760	A
1	AA	1764	A
1	AA	1766	G
1	AA	1767	U
1	AA	1778	G
1	AA	1779	A
1	AA	1780	A
1	AA	1781	C
1	AA	1790	G
1	AA	1791	G
1	AA	1794	C
1	AA	1795	A
19	A7	10	2MG
19	A7	11	C
19	A7	17	H2U
19	A7	20	G
19	A7	23	A
19	A7	44	A
19	A7	47	U
19	A7	76	A
51	B3	2	G
51	B3	7	G
51	B3	10	C
51	B3	11	A
51	B3	13	A
51	B3	14	U

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Mol	Chain	Res	Type
51	B3	24	A
51	B3	34	C
51	B3	36	C
51	B3	40	C
51	B3	41	G
51	B3	42	A
51	B3	43	U
51	B3	44	C
51	B3	45	A
51	B3	46	A
51	B3	48	U
51	B3	50	U
51	B3	56	G
51	B3	59	G
51	B3	61	U
51	B3	70	A
51	B3	71	C
51	B3	74	A
51	B3	81	U
51	B3	82	A
51	B3	84	G
51	B3	86	G
51	B3	88	G
51	B3	89	A
51	B3	90	C
51	B3	91	C
51	B3	92	A
51	B3	95	C
51	B3	98	G
51	B3	99	A
51	B3	101	A
51	B3	102	C
51	B3	105	A
51	B3	109	G
52	B4	4	C
52	B4	11	C
52	B4	12	A
52	B4	14	C
52	B4	18	U
52	B4	20	U
52	B4	22	U
52	B4	23	U

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Mol	Chain	Res	Type
52	B4	24	G
52	B4	34	U
52	B4	48	A
52	B4	55	U
52	B4	58	G
52	B4	59	A
52	B4	60	U
52	B4	61	A
52	B4	62	C
52	B4	63	G
52	B4	64	U
52	B4	68	G
52	B4	69	U
52	B4	70	A
52	B4	71	G
52	B4	72	G
52	B4	76	C
52	B4	88	A
52	B4	89	A
52	B4	92	A
52	B4	95	G
52	B4	99	C
52	B4	100	U
52	B4	102	U
52	B4	103	G
52	B4	104	A
52	B4	105	A
52	B4	106	C
52	B4	108	C
52	B4	109	A
52	B4	111	A
52	B4	114	G
52	B4	115	C
52	B4	118	C
52	B4	128	U
52	B4	133	G
52	B4	134	G
52	B4	135	G
52	B4	136	G
52	B4	137	C
52	B4	138	A
52	B4	139	U

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Mol	Chain	Res	Type
52	B4	140	G
52	B4	145	U
52	B4	147	U
52	B4	153	U
52	B4	156	U
53	B5	4	U
53	B5	6	A
53	B5	7	C
53	B5	13	A
53	B5	15	C
53	B5	16	A
53	B5	17	G
53	B5	19	U
53	B5	33	G
53	B5	39	A
53	B5	40	A
53	B5	43	A
53	B5	44	U
53	B5	46	U
53	B5	60	A
53	B5	62	A
53	B5	66	A
53	B5	68	C
53	B5	71	A
53	B5	74	G
53	B5	77	A
53	B5	85	A
53	B5	89	A
53	B5	92	G
53	B5	99	A
53	B5	108	A
53	B5	109	A
53	B5	110	G
53	B5	112	U
53	B5	114	A
53	B5	115	A
53	B5	118	U
53	B5	120	G
53	B5	125	C
53	B5	128	G
53	B5	129	U
53	B5	131	C

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Mol	Chain	Res	Type
53	B5	132	C
53	B5	133	U
53	B5	134	U
53	B5	135	C
53	B5	149	U
53	B5	150	A
53	B5	151	A
53	B5	160	G
53	B5	161	G
53	B5	163	C
53	B5	170	G
53	B5	171	G
53	B5	175	C
53	B5	186	U
53	B5	189	G
53	B5	199	A
53	B5	200	C
53	B5	202	G
53	B5	210	U
53	B5	211	A
53	B5	214	G
53	B5	216	G
53	B5	217	U
53	B5	218	G
53	B5	219	A
53	B5	220	G
53	B5	230	U
53	B5	240	U
53	B5	250	U
53	B5	251	G
53	B5	252	U
53	B5	265	A
53	B5	268	A
53	B5	269	G
53	B5	272	G
53	B5	274	G
53	B5	278	U
53	B5	279	U
53	B5	280	U
53	B5	281	G
53	B5	287	G
53	B5	288	C

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Mol	Chain	Res	Type
53	B5	291	C
53	B5	298	U
53	B5	304	G
53	B5	310	U
53	B5	315	C
53	B5	318	A
53	B5	330	G
53	B5	338	A
53	B5	341	G
53	B5	343	U
53	B5	352	A
53	B5	354	U
53	B5	355	A
53	B5	356	C
53	B5	362	U
53	B5	364	G
53	B5	368	G
53	B5	370	U
53	B5	372	A
53	B5	373	A
53	B5	375	A
53	B5	376	G
53	B5	384	A
53	B5	387	A
53	B5	391	A
53	B5	394	G
53	B5	398	A
53	B5	399	A
53	B5	402	A
53	B5	403	C
53	B5	404	G
53	B5	406	G
53	B5	411	U
53	B5	413	U
53	B5	414	U
53	B5	415	G
53	B5	423	A
53	B5	434	U
53	B5	438	A
53	B5	440	A
53	B5	441	U
53	B5	451	U

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Mol	Chain	Res	Type
53	B5	453	C
53	B5	455	C
53	B5	457	C
53	B5	465	U
53	B5	467	U
53	B5	471	U
53	B5	476	G
53	B5	477	A
53	B5	478	A
53	B5	479	U
53	B5	480	C
53	B5	481	U
53	B5	482	C
53	B5	483	G
53	B5	484	C
53	B5	485	A
53	B5	486	U
53	B5	492	U
53	B5	494	G
53	B5	496	C
53	B5	500	C
53	B5	513	G
53	B5	515	C
53	B5	532	A
53	B5	533	A
53	B5	547	G
53	B5	548	G
53	B5	553	U
53	B5	554	A
53	B5	555	U
53	B5	556	U
53	B5	570	A
53	B5	572	A
53	B5	576	C
53	B5	578	A
53	B5	585	A
53	B5	590	G
53	B5	592	G
53	B5	596	U
53	B5	597	G
53	B5	602	A
53	B5	604	G

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Mol	Chain	Res	Type
53	B5	605	U
53	B5	606	C
53	B5	607	A
53	B5	608	A
53	B5	609	G
53	B5	611	A
53	B5	612	U
53	B5	616	G
53	B5	619	A
53	B5	620	U
53	B5	621	A
53	B5	622	A
53	B5	627	U
53	B5	629	U
53	B5	636	C
53	B5	638	C
53	B5	639	G
53	B5	647	A
53	B5	648	C
53	B5	649	A
53	B5	652	G
53	B5	653	A
53	B5	654	C
53	B5	663	C
53	B5	664	U
53	B5	677	A
53	B5	678	G
53	B5	681	U
53	B5	688	G
53	B5	689	U
53	B5	690	A
53	B5	691	A
53	B5	695	C
53	B5	704	U
53	B5	705	A
53	B5	706	A
53	B5	708	G
53	B5	710	A
53	B5	711	A
53	B5	712	G
53	B5	713	U
53	B5	715	A

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Mol	Chain	Res	Type
53	B5	716	G
53	B5	717	C
53	B5	718	G
53	B5	735	A
53	B5	736	A
53	B5	742	G
53	B5	743	C
53	B5	744	A
53	B5	748	U
53	B5	766	U
53	B5	767	U
53	B5	783	A
53	B5	784	A
53	B5	785	G
53	B5	786	A
53	B5	787	G
53	B5	800	G
53	B5	801	A
53	B5	802	C
53	B5	808	A
53	B5	817	A
53	B5	819	U
53	B5	820	A
53	B5	833	G
53	B5	837	A
53	B5	844	G
53	B5	846	A
53	B5	848	A
53	B5	851	C
53	B5	856	G
53	B5	858	A
53	B5	860	G
53	B5	861	C
53	B5	879	U
53	B5	880	G
53	B5	882	A
53	B5	884	A
53	B5	894	G
53	B5	896	A
53	B5	907	G
53	B5	913	A
53	B5	914	A

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Mol	Chain	Res	Type
53	B5	915	A
53	B5	916	G
53	B5	917	A
53	B5	921	A
53	B5	923	C
53	B5	925	A
53	B5	932	U
53	B5	933	A
53	B5	934	G
53	B5	937	G
53	B5	944	C
53	B5	951	A
53	B5	959	C
53	B5	962	A
53	B5	963	G
53	B5	965	A
53	B5	970	A
53	B5	971	C
53	B5	972	G
53	B5	973	A
53	B5	978	C
53	B5	979	G
53	B5	987	U
53	B5	990	A
53	B5	994	G
53	B5	1009	A
53	B5	1011	A
53	B5	1017	C
53	B5	1024	G
53	B5	1025	A
53	B5	1027	A
53	B5	1029	G
53	B5	1030	A
53	B5	1033	U
53	B5	1034	U
53	B5	1035	G
53	B5	1036	A
53	B5	1041	U
53	B5	1047	A
53	B5	1054	A
53	B5	1056	U
53	B5	1065	A

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Mol	Chain	Res	Type
53	B5	1066	G
53	B5	1078	U
53	B5	1079	A
53	B5	1080	A
53	B5	1085	A
53	B5	1088	U
53	B5	1093	A
53	B5	1099	A
53	B5	1101	G
53	B5	1104	G
53	B5	1105	A
53	B5	1110	U
53	B5	1111	U
53	B5	1112	A
53	B5	1116	G
53	B5	1117	G
53	B5	1126	G
53	B5	1127	G
53	B5	1128	U
53	B5	1129	A
53	B5	1133	A
53	B5	1135	A
53	B5	1142	G
53	B5	1144	U
53	B5	1150	A
53	B5	1153	A
53	B5	1160	C
53	B5	1162	U
53	B5	1164	G
53	B5	1166	G
53	B5	1173	U
53	B5	1174	G
53	B5	1175	C
53	B5	1177	G
53	B5	1181	U
53	B5	1182	A
53	B5	1191	U
53	B5	1200	A
53	B5	1201	C
53	B5	1208	U
53	B5	1219	C
53	B5	1227	C

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Mol	Chain	Res	Type
53	B5	1228	C
53	B5	1229	G
53	B5	1231	A
53	B5	1232	C
53	B5	1235	U
53	B5	1236	G
53	B5	1237	G
53	B5	1238	C
53	B5	1244	A
53	B5	1245	A
53	B5	1248	C
53	B5	1250	G
53	B5	1251	A
53	B5	1253	U
53	B5	1254	C
53	B5	1257	C
53	B5	1258	U
53	B5	1263	A
53	B5	1264	G
53	B5	1268	G
53	B5	1273	A
53	B5	1274	A
53	B5	1279	C
53	B5	1282	G
53	B5	1287	A
53	B5	1291	A
53	B5	1298	C
53	B5	1303	A
53	B5	1304	A
53	B5	1305	U
53	B5	1306	G
53	B5	1307	G
53	B5	1308	A
53	B5	1309	U
53	B5	1311	G
53	B5	1313	G
53	B5	1318	A
53	B5	1319	G
53	B5	1347	U
53	B5	1357	G
53	B5	1358	C
53	B5	1359	C

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Mol	Chain	Res	Type
53	B5	1360	C
53	B5	1361	U
53	B5	1363	A
53	B5	1370	G
53	B5	1384	U
53	B5	1385	C
53	B5	1391	C
53	B5	1392	G
53	B5	1402	C
53	B5	1404	G
53	B5	1406	A
53	B5	1407	A
53	B5	1408	G
53	B5	1412	G
53	B5	1418	A
53	B5	1421	G
53	B5	1422	G
53	B5	1423	C
53	B5	1425	U
53	B5	1426	C
53	B5	1429	G
53	B5	1431	G
53	B5	1434	G
53	B5	1439	U
53	B5	1442	U
53	B5	1447	G
53	B5	1452	A
53	B5	1453	A
53	B5	1455	U
53	B5	1456	A
53	B5	1457	U
53	B5	1481	A
53	B5	1482	A
53	B5	1507	G
53	B5	1510	G
53	B5	1515	A
53	B5	1526	U
53	B5	1529	A
53	B5	1530	U
53	B5	1532	C
53	B5	1533	U
53	B5	1548	C

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Mol	Chain	Res	Type
53	B5	1550	C
53	B5	1554	U
53	B5	1555	U
53	B5	1571	A
53	B5	1572	U
53	B5	1580	A
53	B5	1582	C
53	B5	1583	A
53	B5	1584	U
53	B5	1588	A
53	B5	1590	G
53	B5	1591	G
53	B5	1592	G
53	B5	1593	A
53	B5	1598	G
53	B5	1602	A
53	B5	1607	U
53	B5	1617	G
53	B5	1622	U
53	B5	1623	G
53	B5	1624	G
53	B5	1627	U
53	B5	1628	C
53	B5	1637	A
53	B5	1639	C
53	B5	1640	G
53	B5	1642	A
53	B5	1644	C
53	B5	1647	A
53	B5	1648	A
53	B5	1649	U
53	B5	1650	G
53	B5	1657	C
53	B5	1658	G
53	B5	1666	G
53	B5	1668	G
53	B5	1669	C
53	B5	1670	C
53	B5	1671	C
53	B5	1672	U
53	B5	1681	U
53	B5	1682	U

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Mol	Chain	Res	Type
53	B5	1683	A
53	B5	1688	U
53	B5	1689	U
53	B5	1697	A
53	B5	1698	C
53	B5	1701	C
53	B5	1702	U
53	B5	1703	U
53	B5	1704	A
53	B5	1709	C
53	B5	1710	C
53	B5	1712	G
53	B5	1715	A
53	B5	1718	G
53	B5	1719	G
53	B5	1720	U
53	B5	1721	U
53	B5	1722	U
53	B5	1727	G
53	B5	1729	A
53	B5	1733	G
53	B5	1735	G
53	B5	1736	G
53	B5	1737	U
53	B5	1738	C
53	B5	1744	G
53	B5	1745	C
53	B5	1746	U
53	B5	1750	A
53	B5	1751	G
53	B5	1759	C
53	B5	1762	C
53	B5	1764	U
53	B5	1765	U
53	B5	1768	U
53	B5	1769	G
53	B5	1774	C
53	B5	1775	G
53	B5	1777	U
53	B5	1784	G
53	B5	1797	A
53	B5	1805	C

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Mol	Chain	Res	Type
53	B5	1806	A
53	B5	1813	A
53	B5	1814	A
53	B5	1817	G
53	B5	1819	U
53	B5	1824	U
53	B5	1826	C
53	B5	1830	G
53	B5	1832	C
53	B5	1840	U
53	B5	1841	A
53	B5	1842	A
53	B5	1849	C
53	B5	1853	U
53	B5	1855	U
53	B5	1859	A
53	B5	1864	A
53	B5	1866	C
53	B5	1867	A
53	B5	1880	U
53	B5	1886	A
53	B5	1889	G
53	B5	1890	U
53	B5	1892	G
53	B5	1895	A
53	B5	1900	A
53	B5	1901	A
53	B5	1902	G
53	B5	1906	G
53	B5	1909	A
53	B5	1918	C
53	B5	1919	G
53	B5	1927	G
53	B5	1932	A
53	B5	1933	A
53	B5	1934	G
53	B5	1935	G
53	B5	1946	A
53	B5	2102	U
53	B5	2103	U
53	B5	2104	A
53	B5	2114	C

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Mol	Chain	Res	Type
53	B5	2115	G
53	B5	2116	G
53	B5	2120	A
53	B5	2122	G
53	B5	2125	A
53	B5	2126	A
53	B5	2131	A
53	B5	2134	G
53	B5	2138	A
53	B5	2139	A
53	B5	2140	U
53	B5	2142	A
53	B5	2159	U
53	B5	2160	G
53	B5	2161	G
53	B5	2167	A
53	B5	2170	U
53	B5	2173	U
53	B5	2174	G
53	B5	2175	U
53	B5	2178	A
53	B5	2189	U
53	B5	2199	G
53	B5	2208	A
53	B5	2209	U
53	B5	2210	G
53	B5	2212	C
53	B5	2213	A
53	B5	2214	A
53	B5	2215	A
53	B5	2216	G
53	B5	2217	U
53	B5	2219	A
53	B5	2220	A
53	B5	2221	G
53	B5	2222	A
53	B5	2223	A
53	B5	2224	A
53	B5	2225	U
53	B5	2226	U
53	B5	2227	C
53	B5	2228	A

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Mol	Chain	Res	Type
53	B5	2229	A
53	B5	2230	C
53	B5	2237	C
53	B5	2242	A
53	B5	2244	A
53	B5	2249	G
53	B5	2252	A
53	B5	2255	A
53	B5	2256	A
53	B5	2257	C
53	B5	2259	A
53	B5	2263	C
53	B5	2273	G
53	B5	2279	A
53	B5	2285	C
53	B5	2286	U
53	B5	2287	C
53	B5	2288	G
53	B5	2297	U
53	B5	2298	U
53	B5	2301	U
53	B5	2303	A
53	B5	2306	C
53	B5	2308	C
53	B5	2309	A
53	B5	2310	U
53	B5	2313	A
53	B5	2314	U
53	B5	2315	G
53	B5	2317	A
53	B5	2319	U
53	B5	2326	A
53	B5	2335	G
53	B5	2336	U
53	B5	2340	U
53	B5	2347	U
53	B5	2350	C
53	B5	2355	G
53	B5	2360	C
53	B5	2362	C
53	B5	2363	A
53	B5	2365	C

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Mol	Chain	Res	Type
53	B5	2368	A
53	B5	2371	G
53	B5	2372	A
53	B5	2373	A
53	B5	2374	C
53	B5	2379	U
53	B5	2383	C
53	B5	2385	G
53	B5	2388	U
53	B5	2394	G
53	B5	2397	A
53	B5	2402	A
53	B5	2403	G
53	B5	2411	U
53	B5	2415	C
53	B5	2417	U
53	B5	2419	A
53	B5	2435	G
53	B5	2446	U
53	B5	2452	G
53	B5	2454	G
53	B5	2457	G
53	B5	2458	A
53	B5	2460	U
53	B5	2461	A
53	B5	2462	A
53	B5	2463	G
53	B5	2464	U
53	B5	2468	A
53	B5	2469	G
53	B5	2470	C
53	B5	2472	U
53	B5	2473	C
53	B5	2474	G
53	B5	2475	G
53	B5	2477	G
53	B5	2478	C
53	B5	2479	C
53	B5	2484	A
53	B5	2487	U
53	B5	2494	A
53	B5	2497	U

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Mol	Chain	Res	Type
53	B5	2498	U
53	B5	2499	U
53	B5	2506	U
53	B5	2510	U
53	B5	2514	U
53	B5	2515	A
53	B5	2521	U
53	B5	2524	A
53	B5	2539	A
53	B5	2540	A
53	B5	2541	U
53	B5	2542	U
53	B5	2543	U
53	B5	2546	C
53	B5	2548	C
53	B5	2550	U
53	B5	2551	C
53	B5	2552	C
53	B5	2556	U
53	B5	2560	U
53	B5	2561	A
53	B5	2562	G
53	B5	2570	U
53	B5	2572	C
53	B5	2573	G
53	B5	2575	G
53	B5	2576	G
53	B5	2577	C
53	B5	2578	U
53	B5	2585	G
53	B5	2586	G
53	B5	2593	A
53	B5	2596	U
53	B5	2606	G
53	B5	2607	G
53	B5	2608	G
53	B5	2610	G
53	B5	2618	G
53	B5	2619	G
53	B5	2623	G
53	B5	2638	C
53	B5	2652	U

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Mol	Chain	Res	Type
53	B5	2655	U
53	B5	2656	A
53	B5	2658	G
53	B5	2674	A
53	B5	2675	C
53	B5	2676	A
53	B5	2677	G
53	B5	2678	A
53	B5	2680	A
53	B5	2690	G
53	B5	2691	A
53	B5	2694	A
53	B5	2698	G
53	B5	2703	A
53	B5	2704	A
53	B5	2706	G
53	B5	2714	G
53	B5	2715	A
53	B5	2718	U
53	B5	2719	U
53	B5	2725	U
53	B5	2735	U
53	B5	2736	A
53	B5	2749	G
53	B5	2752	U
53	B5	2755	C
53	B5	2757	U
53	B5	2758	A
53	B5	2765	C
53	B5	2773	C
53	B5	2774	C
53	B5	2775	U
53	B5	2776	C
53	B5	2778	G
53	B5	2780	A
53	B5	2781	U
53	B5	2783	U
53	B5	2788	C
53	B5	2794	G
53	B5	2795	U
53	B5	2796	G
53	B5	2797	C

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Mol	Chain	Res	Type
53	B5	2799	A
53	B5	2800	G
53	B5	2801	A
53	B5	2802	A
53	B5	2809	C
53	B5	2810	C
53	B5	2816	G
53	B5	2817	A
53	B5	2818	U
53	B5	2827	U
53	B5	2829	U
53	B5	2830	G
53	B5	2837	A
53	B5	2838	A
53	B5	2839	G
53	B5	2840	C
53	B5	2841	G
53	B5	2842	U
53	B5	2845	A
53	B5	2849	C
53	B5	2850	G
53	B5	2852	C
53	B5	2853	A
53	B5	2860	U
53	B5	2863	G
53	B5	2864	A
53	B5	2865	U
53	B5	2866	U
53	B5	2867	C
53	B5	2868	U
53	B5	2872	A
53	B5	2874	G
53	B5	2875	U
53	B5	2877	G
53	B5	2882	U
53	B5	2887	A
53	B5	2888	U
53	B5	2899	C
53	B5	2902	A
53	B5	2904	U
53	B5	2905	U
53	B5	2910	A

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Mol	Chain	Res	Type
53	B5	2914	G
53	B5	2915	U
53	B5	2919	A
53	B5	2921	U
53	B5	2931	C
53	B5	2936	A
53	B5	2941	A
53	B5	2942	C
53	B5	2943	G
53	B5	2944	U
53	B5	2945	G
53	B5	2947	G
53	B5	2953	U
53	B5	2962	U
53	B5	2965	U
53	B5	2975	U
53	B5	2978	U
53	B5	2980	U
53	B5	2981	U
53	B5	2988	C
53	B5	2989	U
53	B5	2996	U
53	B5	2997	G
53	B5	3001	C
53	B5	3002	C
53	B5	3004	C
53	B5	3010	U
53	B5	3012	A
53	B5	3019	U
53	B5	3028	G
53	B5	3030	G
53	B5	3034	C
53	B5	3040	A
53	B5	3046	A
53	B5	3048	A
53	B5	3049	A
53	B5	3050	U
53	B5	3051	U
53	B5	3052	G
53	B5	3055	U
53	B5	3056	U
53	B5	3057	U

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Mol	Chain	Res	Type
53	B5	3058	U
53	B5	3061	G
53	B5	3062	G
53	B5	3081	C
53	B5	3083	G
53	B5	3084	C
53	B5	3086	A
53	B5	3088	G
53	B5	3100	U
53	B5	3112	G
53	B5	3113	A
53	B5	3114	A
53	B5	3116	G
53	B5	3120	C
53	B5	3121	U
53	B5	3122	A
53	B5	3123	A
53	B5	3124	G
53	B5	3126	C
53	B5	3134	A
53	B5	3136	G
53	B5	3139	A
53	B5	3140	G
53	B5	3141	A
53	B5	3142	A
53	B5	3143	C
53	B5	3144	G
53	B5	3149	G
53	B5	3151	U
53	B5	3152	U
53	B5	3154	C
53	B5	3158	G
53	B5	3159	C
53	B5	3160	U
53	B5	3161	C
53	B5	3162	C
53	B5	3165	A
53	B5	3166	C
53	B5	3167	A
53	B5	3170	A
53	B5	3171	U
53	B5	3174	A

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Mol	Chain	Res	Type
53	B5	3175	C
53	B5	3176	C
53	B5	3182	U
53	B5	3183	A
53	B5	3184	U
53	B5	3186	A
53	B5	3188	G
53	B5	3198	U
53	B5	3199	G
53	B5	3200	G
53	B5	3205	U
53	B5	3206	A
53	B5	3208	C
53	B5	3211	U
53	B5	3216	U
53	B5	3217	A
53	B5	3220	G
53	B5	3229	G
53	B5	3230	G
53	B5	3235	C
53	B5	3243	A
53	B5	3245	A
53	B5	3246	G
53	B5	3247	G
53	B5	3248	C
53	B5	3249	C
53	B5	3252	G
53	B5	3253	G
53	B5	3259	U
53	B5	3266	G
53	B5	3267	U
53	B5	3277	C
53	B5	3290	G
53	B5	3295	A
53	B5	3296	A
53	B5	3298	C
53	B5	3304	U
53	B5	3305	A
53	B5	3306	U
53	B5	3307	A
53	B5	3317	U
53	B5	3320	A

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Mol	Chain	Res	Type
53	B5	3325	G
53	B5	3326	G
53	B5	3327	G
53	B5	3328	G
53	B5	3330	A
53	B5	3331	U
53	B5	3332	U
53	B5	3333	G
53	B5	3334	U
53	B5	3337	G
53	B5	3338	C
53	B5	3354	U
53	B5	3355	U
53	B5	3364	C
53	B5	3368	U
53	B5	3369	G
53	B5	3374	U
53	B5	3385	U
53	B5	3386	G
53	B5	3389	U
53	B5	3391	A
53	B5	3395	G
53	B5	3396	U

All (235) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
1	AA	2	A
1	AA	42	G
1	AA	47	A
1	AA	56	U
1	AA	60	U
1	AA	66	U
1	AA	109	G
1	AA	175	G
1	AA	199	G
1	AA	212	U
1	AA	229	U
1	AA	312	A
1	AA	350	U
1	AA	365	G
1	AA	400	A

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Mol	Chain	Res	Type
1	AA	438	A
1	AA	444	C
1	AA	461	G
1	AA	478	A
1	AA	502	U
1	AA	578	U
1	AA	652	G
1	AA	748	U
1	AA	759	U
1	AA	770	A
1	AA	796	A
1	AA	852	C
1	AA	857	U
1	AA	864	U
1	AA	922	A
1	AA	1049	U
1	AA	1055	U
1	AA	1090	A
1	AA	1098	G
1	AA	1133	U
1	AA	1134	A
1	AA	1159	U
1	AA	1181	U
1	AA	1192	A
1	AA	1196	G
1	AA	1199	A
1	AA	1216	C
1	AA	1240	A
1	AA	1241	G
1	AA	1265	U
1	AA	1281	C
1	AA	1318	A
1	AA	1328	A
1	AA	1341	A
1	AA	1378	U
1	AA	1429	U
1	AA	1457	C
1	AA	1478	G
1	AA	1488	C
1	AA	1520	U
1	AA	1521	G
1	AA	1535	C

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Mol	Chain	Res	Type
1	AA	1545	A
1	AA	1546	G
1	AA	1547	C
1	AA	1560	G
1	AA	1566	C
1	AA	1579	C
1	AA	1583	U
1	AA	1584	A
1	AA	1612	A
1	AA	1632	C
1	AA	1634	C
1	AA	1635	C
1	AA	1636	G
1	AA	1639	C
1	AA	1654	U
1	AA	1655	U
1	AA	1675	C
1	AA	1759	U
1	AA	1766	G
1	AA	1779	A
1	AA	1789	A
1	AA	1790	G
19	A7	10	2MG
19	A7	22	G
19	A7	41	U
51	B3	23	A
51	B3	70	A
51	B3	89	A
51	B3	91	C
52	B4	19	C
52	B4	21	C
52	B4	57	C
52	B4	60	U
52	B4	61	A
52	B4	69	U
52	B4	71	G
52	B4	102	U
52	B4	106	C
52	B4	152	G
53	B5	114	A
53	B5	119	U
53	B5	133	U

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Mol	Chain	Res	Type
53	B5	134	U
53	B5	210	U
53	B5	215	G
53	B5	239	G
53	B5	250	U
53	B5	251	G
53	B5	267	G
53	B5	277	G
53	B5	279	U
53	B5	355	A
53	B5	372	A
53	B5	374	A
53	B5	397	A
53	B5	475	G
53	B5	477	A
53	B5	481	U
53	B5	554	A
53	B5	596	U
53	B5	620	U
53	B5	621	A
53	B5	638	C
53	B5	677	A
53	B5	705	A
53	B5	711	A
53	B5	712	G
53	B5	714	G
53	B5	715	A
53	B5	735	A
53	B5	742	G
53	B5	786	A
53	B5	801	A
53	B5	860	G
53	B5	914	A
53	B5	932	U
53	B5	970	A
53	B5	977	U
53	B5	979	G
53	B5	1033	U
53	B5	1079	A
53	B5	1084	A
53	B5	1163	A
53	B5	1181	U

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Mol	Chain	Res	Type
53	B5	1200	A
53	B5	1235	U
53	B5	1240	A
53	B5	1244	A
53	B5	1258	U
53	B5	1263	A
53	B5	1302	A
53	B5	1303	A
53	B5	1318	A
53	B5	1331	U
53	B5	1347	U
53	B5	1390	A
53	B5	1514	G
53	B5	1549	U
53	B5	1553	U
53	B5	1582	C
53	B5	1583	A
53	B5	1590	G
53	B5	1623	G
53	B5	1626	U
53	B5	1643	A
53	B5	1648	A
53	B5	1654	A
53	B5	1666	G
53	B5	1720	U
53	B5	1721	U
53	B5	1737	U
53	B5	1776	G
53	B5	1818	U
53	B5	1842	A
53	B5	1900	A
53	B5	1925	U
53	B5	1947	G
53	B5	2138	A
53	B5	2174	G
53	B5	2177	G
53	B5	2214	A
53	B5	2219	A
53	B5	2224	A
53	B5	2225	U
53	B5	2229	A
53	B5	2283	G

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Mol	Chain	Res	Type
53	B5	2286	U
53	B5	2297	U
53	B5	2314	U
53	B5	2360	C
53	B5	2373	A
53	B5	2459	A
53	B5	2468	A
53	B5	2469	G
53	B5	2474	G
53	B5	2477	G
53	B5	2520	A
53	B5	2538	U
53	B5	2569	A
53	B5	2571	U
53	B5	2618	G
53	B5	2644	C
53	B5	2651	G
53	B5	2654	C
53	B5	2677	G
53	B5	2705	A
53	B5	2714	G
53	B5	2778	G
53	B5	2787	G
53	B5	2796	G
53	B5	2830	G
53	B5	2837	A
53	B5	2875	U
53	B5	2913	C
53	B5	2941	A
53	B5	2988	C
53	B5	3039	C
53	B5	3045	G
53	B5	3049	A
53	B5	3051	U
53	B5	3077	A
53	B5	3083	G
53	B5	3110	C
53	B5	3113	A
53	B5	3121	U
53	B5	3138	U
53	B5	3141	A
53	B5	3142	A

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Mol	Chain	Res	Type
53	B5	3143	C
53	B5	3160	U
53	B5	3195	U
53	B5	3219	U
53	B5	3246	G
53	B5	3304	U
53	B5	3313	U
53	B5	3330	A
53	B5	3384	U
53	B5	3385	U

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

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

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
19	2MG	A7	10	19	18,26,27	1.88	8 (44%)	21,38,41	2.83	10 (47%)
19	H2U	A7	16	19	17,21,22	1.89	3 (17%)	23,30,33	2.37	9 (39%)
19	H2U	A7	17	19	17,21,22	2.68	7 (41%)	23,30,33	2.71	6 (26%)
19	M2G	A7	26	19	18,27,28	1.81	4 (22%)	22,40,43	2.92	9 (40%)
19	OMC	A7	32	19	15,22,23	1.73	3 (20%)	20,31,34	2.07	5 (25%)
19	OMG	A7	34	19	18,26,27	1.67	3 (16%)	21,38,41	2.98	3 (14%)
19	YYG	A7	37	19	28,42,43	1.74	7 (25%)	28,62,65	2.82	13 (46%)
19	PSU	A7	39	19	15,21,22	1.84	7 (46%)	16,30,33	2.59	4 (25%)
19	5MC	A7	40	19	14,22,23	2.38	4 (28%)	17,32,35	2.07	4 (23%)
19	7MG	A7	46	19	20,26,27	1.68	5 (25%)	23,39,42	2.56	7 (30%)
19	5MC	A7	49	19	14,22,23	1.77	2 (14%)	17,32,35	2.25	5 (29%)
19	5MU	A7	54	19	13,22,23	0.96	0	16,32,35	4.68	6 (37%)
19	PSU	A7	55	19	15,21,22	1.65	2 (13%)	16,30,33	3.69	4 (25%)
19	1MA	A7	58	19	15,25,26	0.91	1 (6%)	15,37,40	2.61	4 (26%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
19	2MG	A7	10	19	-	0/5/27/28	0/3/3/3
19	H2U	A7	16	19	-	0/7/38/39	0/2/2/2
19	H2U	A7	17	19	-	0/7/38/39	0/2/2/2
19	M2G	A7	26	19	-	0/7/29/30	0/3/3/3
19	OMC	A7	32	19	-	0/5/27/28	0/2/2/2
19	OMG	A7	34	19	-	0/5/27/28	0/3/3/3
19	YYG	A7	37	19	-	0/20/42/43	0/4/4/4
19	PSU	A7	39	19	-	0/7/25/26	0/2/2/2
19	5MC	A7	40	19	-	0/3/25/26	0/2/2/2
19	7MG	A7	46	19	-	0/7/37/38	0/3/3/3
19	5MC	A7	49	19	-	0/3/25/26	0/2/2/2
19	5MU	A7	54	19	-	0/3/25/26	0/2/2/2
19	PSU	A7	55	19	-	0/7/25/26	0/2/2/2
19	1MA	A7	58	19	-	0/3/25/26	0/3/3/3

All (56) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
19	A7	17	H2U	C4-N3	-6.07	1.28	1.37
19	A7	17	H2U	C2-N3	-5.93	1.26	1.38
19	A7	55	PSU	C6-C5	-4.48	1.32	1.38
19	A7	40	5MC	C2'-C1'	-4.37	1.46	1.53
19	A7	40	5MC	C4-N3	-4.26	1.28	1.35
19	A7	37	YYG	C3-N3	-3.91	1.43	1.49
19	A7	16	H2U	C2-N3	-3.66	1.31	1.38
19	A7	16	H2U	C4-N3	-3.63	1.31	1.37
19	A7	26	M2G	O4'-C1'	-3.58	1.36	1.41
19	A7	10	2MG	C2-N2	-3.37	1.30	1.34
19	A7	34	OMG	O4'-C1'	-3.20	1.36	1.41
19	A7	37	YYG	C4-N3	-3.09	1.35	1.39
19	A7	37	YYG	C2'-C1'	-2.97	1.48	1.53
19	A7	32	OMC	O2'-C2'	-2.94	1.34	1.42
19	A7	39	PSU	O2'-C2'	-2.93	1.36	1.43
19	A7	10	2MG	O2'-C2'	-2.88	1.36	1.43
19	A7	39	PSU	O4'-C1'	-2.76	1.40	1.44
19	A7	26	M2G	C8-N7	-2.68	1.29	1.34
19	A7	46	7MG	C8-N9	-2.62	1.41	1.45
19	A7	39	PSU	C5-C1'	-2.62	1.50	1.52
19	A7	17	H2U	O4'-C4'	-2.53	1.39	1.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
19	A7	39	PSU	C2'-C1'	-2.39	1.51	1.53
19	A7	39	PSU	O4'-C4'	-2.32	1.39	1.45
19	A7	10	2MG	C4-N3	-2.23	1.32	1.35
19	A7	32	OMC	C2-N3	-2.22	1.33	1.38
19	A7	10	2MG	C5-C4	-2.13	1.35	1.40
19	A7	17	H2U	C2'-C1'	-2.09	1.46	1.53
19	A7	46	7MG	C4-N3	-2.04	1.31	1.34
19	A7	39	PSU	C6-C5	-2.00	1.35	1.38
19	A7	39	PSU	C2-N1	-2.00	1.34	1.38
19	A7	37	YYG	C6-C5	2.04	1.44	1.40
19	A7	10	2MG	C6-N1	2.16	1.36	1.33
19	A7	55	PSU	C2'-C1'	2.20	1.55	1.53
19	A7	37	YYG	C2-N2	2.22	1.38	1.35
19	A7	37	YYG	O4'-C1'	2.23	1.44	1.41
19	A7	10	2MG	O4'-C1'	2.34	1.44	1.41
19	A7	58	1MA	C6-N6	2.36	1.33	1.29
19	A7	26	M2G	C5'-C4'	2.37	1.59	1.51
19	A7	10	2MG	C6-C5	2.39	1.46	1.41
19	A7	37	YYG	C14-C15	2.41	1.59	1.53
19	A7	17	H2U	C2-N1	2.46	1.39	1.35
19	A7	10	2MG	C2'-C3'	2.62	1.60	1.53
19	A7	46	7MG	CM7-N7	2.74	1.50	1.46
19	A7	46	7MG	C6-C5	2.80	1.45	1.41
19	A7	34	OMG	C6-N1	3.11	1.38	1.33
19	A7	17	H2U	C6-N1	3.42	1.51	1.47
19	A7	17	H2U	C1'-N1	3.49	1.53	1.45
19	A7	49	5MC	CM5-C5	3.55	1.58	1.51
19	A7	46	7MG	C6-N1	3.55	1.39	1.33
19	A7	34	OMG	C3'-C2'	3.73	1.61	1.53
19	A7	40	5MC	C5-C4	3.83	1.47	1.41
19	A7	32	OMC	O4'-C1'	3.86	1.46	1.41
19	A7	26	M2G	C6-N1	4.24	1.40	1.33
19	A7	40	5MC	O4'-C1'	4.24	1.47	1.41
19	A7	16	H2U	C2-N1	4.26	1.42	1.35
19	A7	49	5MC	C5-C4	4.32	1.48	1.41

All (89) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
19	A7	54	5MU	C5-C4-N3	-12.29	115.03	125.35
19	A7	34	OMG	C5-C6-N1	-10.12	110.29	123.52
19	A7	26	M2G	C5-C6-N1	-8.97	111.80	123.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
19	A7	46	7MG	C5-C6-N1	-7.67	111.97	123.39
19	A7	58	1MA	C1'-N9-C4	-7.61	118.31	126.81
19	A7	54	5MU	C4'-O4'-C1'	-6.17	103.10	109.64
19	A7	40	5MC	C4'-O4'-C1'	-6.00	103.28	109.64
19	A7	10	2MG	C6-C5-C4	-5.75	114.29	120.86
19	A7	16	H2U	C6-N1-C2	-5.43	113.77	122.16
19	A7	55	PSU	O2'-C2'-C1'	-5.18	100.66	111.93
19	A7	10	2MG	C5-C6-N1	-4.66	117.43	123.52
19	A7	39	PSU	C4'-O4'-C1'	-4.10	105.31	109.54
19	A7	37	YYG	O23-C21-O22	-4.00	118.99	124.61
19	A7	17	H2U	O2-C2-N1	-3.94	118.01	123.17
19	A7	39	PSU	C5-C6-N1	-3.89	118.96	124.38
19	A7	54	5MU	C5M-C5-C4	-3.87	115.68	119.97
19	A7	40	5MC	N4-C4-N3	-3.74	111.44	116.92
19	A7	26	M2G	N3-C2-N1	-3.74	120.00	126.35
19	A7	17	H2U	C6-N1-C2	-3.69	116.45	122.16
19	A7	49	5MC	O2'-C2'-C1'	-3.52	100.61	111.61
19	A7	26	M2G	C6-C5-C4	-3.46	116.90	120.86
19	A7	46	7MG	C4-N9-C1'	-3.33	118.77	126.65
19	A7	32	OMC	C6-N1-C2	-3.17	116.15	121.33
19	A7	10	2MG	N3-C2-N1	-3.11	121.53	126.19
19	A7	49	5MC	C2'-C3'-C4'	-3.11	96.27	102.64
19	A7	17	H2U	O2-C2-N3	-2.90	115.75	121.44
19	A7	16	H2U	O2'-C2'-C1'	-2.87	100.00	109.98
19	A7	26	M2G	C2-N3-C4	-2.85	111.87	114.99
19	A7	37	YYG	O2'-C2'-C1'	-2.60	103.48	111.61
19	A7	17	H2U	C4-N3-C2	-2.60	123.42	125.77
19	A7	32	OMC	N4-C4-N3	-2.57	112.02	116.50
19	A7	46	7MG	C5-C4-N3	-2.46	124.23	126.74
19	A7	40	5MC	O3'-C3'-C2'	-2.38	104.18	111.86
19	A7	40	5MC	C2'-C3'-C4'	-2.31	97.91	102.64
19	A7	55	PSU	C2'-C3'-C4'	-2.31	97.92	102.64
19	A7	46	7MG	N2-C2-N3	-2.30	113.41	117.20
19	A7	17	H2U	C3'-C2'-C1'	-2.29	96.82	101.44
19	A7	46	7MG	N2-C2-N1	-2.28	113.44	117.20
19	A7	37	YYG	C14-C15-N20	-2.24	106.58	110.81
19	A7	34	OMG	C2'-C3'-C4'	-2.22	96.59	101.85
19	A7	26	M2G	O3'-C3'-C4'	-2.21	104.41	111.01
19	A7	10	2MG	O2'-C2'-C1'	-2.14	104.91	111.61
19	A7	32	OMC	CM2-O2'-C2'	-2.14	108.59	114.58
19	A7	37	YYG	O17-C16-C15	-2.12	118.19	124.08
19	A7	16	H2U	O2-C2-N1	-2.03	120.51	123.17

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
19	A7	10	2MG	C5'-C4'-C3'	-2.02	107.39	115.20
19	A7	16	H2U	O2-C2-N3	-2.01	117.49	121.44
19	A7	55	PSU	C5-C1'-C2'	2.04	118.90	115.44
19	A7	37	YYG	O3'-C3'-C4'	2.10	117.27	111.01
19	A7	26	M2G	O4'-C1'-N9	2.20	112.26	108.11
19	A7	16	H2U	O4'-C4'-C3'	2.22	109.67	105.16
19	A7	54	5MU	O4'-C4'-C5'	2.26	117.38	109.29
19	A7	10	2MG	C4'-O4'-C1'	2.31	112.09	109.64
19	A7	32	OMC	O4'-C1'-N1	2.33	112.53	108.10
19	A7	49	5MC	O5'-C5'-C4'	2.39	117.65	109.09
19	A7	54	5MU	O4'-C4'-C3'	2.44	110.10	105.16
19	A7	58	1MA	O2'-C2'-C1'	2.55	119.59	111.61
19	A7	49	5MC	O3'-C3'-C4'	2.65	118.93	111.01
19	A7	10	2MG	C2-N3-C4	2.67	117.92	114.99
19	A7	16	H2U	C5-C4-N3	2.70	119.47	116.62
19	A7	58	1MA	O4'-C1'-N9	2.92	113.62	108.11
19	A7	37	YYG	C19-O18-C16	2.92	122.90	115.97
19	A7	16	H2U	C5-C6-N1	3.01	114.06	110.76
19	A7	26	M2G	CM1-N2-C2	3.06	124.41	121.34
19	A7	37	YYG	C4'-O4'-C1'	3.06	112.89	109.64
19	A7	37	YYG	O23-C21-N20	3.48	117.88	110.84
19	A7	26	M2G	N1-C2-N2	3.73	121.20	117.14
19	A7	10	2MG	N2-C2-N3	3.79	121.34	116.94
19	A7	16	H2U	C1'-N1-C2	3.96	123.74	118.19
19	A7	58	1MA	C6-C5-C4	4.18	120.16	116.80
19	A7	10	2MG	C6-N1-C2	4.36	121.49	115.24
19	A7	39	PSU	O4'-C1'-C2'	4.47	109.52	104.69
19	A7	37	YYG	C15-N20-C21	4.47	133.14	121.06
19	A7	26	M2G	CM2-N2-C2	4.63	125.99	121.34
19	A7	46	7MG	N1-C2-N3	4.65	133.12	125.51
19	A7	37	YYG	C6-C5-C4	4.70	123.28	119.93
19	A7	46	7MG	C8-N9-C1'	4.95	137.30	122.43
19	A7	37	YYG	C14-C15-C16	5.26	122.53	109.95
19	A7	37	YYG	C3-N3-C4	5.63	126.91	118.41
19	A7	16	H2U	N3-C2-N1	5.77	121.98	116.64
19	A7	10	2MG	CM2-N2-C2	6.13	129.93	123.03
19	A7	37	YYG	C24-O23-C21	6.31	123.52	115.65
19	A7	32	OMC	C6-C5-C4	6.49	119.98	117.44
19	A7	39	PSU	C4-N3-C2	6.54	120.61	115.16
19	A7	49	5MC	C4'-O4'-C1'	6.55	116.58	109.64
19	A7	34	OMG	C6-N1-C2	7.70	124.91	115.88
19	A7	17	H2U	N3-C2-N1	10.13	126.03	116.64

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
19	A7	54	5MU	C4-N3-C2	11.14	124.45	115.16
19	A7	55	PSU	C4-N3-C2	12.99	126.00	115.16

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
19	A7	10	2MG	1	0
19	A7	37	YYG	1	0
19	A7	40	5MC	3	0

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
7	AH	1
18	AT	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	AH	32:LYS	C	33:VAL	N	2.59
1	AT	78:LYS	C	79:LEU	N	2.49