



Full wwPDB EM Validation Report ⓘ

Jun 17, 2026 – 06:47 pm BST

PDB ID : 7NQL / pdb_00007nql
EMDB ID : EMD-12529
Title : 55S mammalian mitochondrial ribosome with ICT1 and P site tRNAMet
Authors : Kummer, E.; Schubert, K.; Ban, N.
Deposited on : 2021-03-01
Resolution : 3.40 Å(reported)

This is a Full wwPDB EM Validation Report for a publicly released PDB entry.

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

A user guide is available at

<https://www.wwpdb.org/validation/2017/EMValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

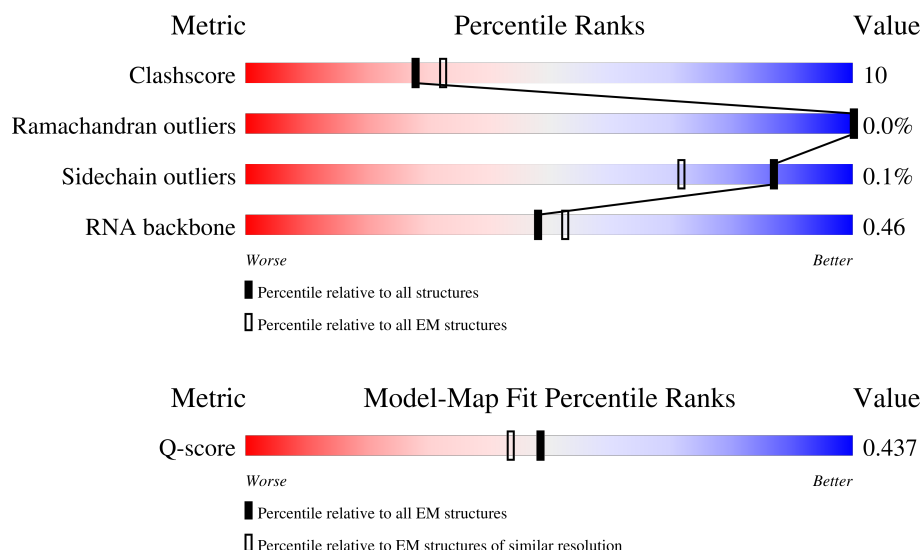
EMDB validation analysis : 0.0.1.dev132
Mogul : 1.8.4, CSD as541be (2020)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

1 Overall quality at a glance

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

The reported resolution of this entry is 3.40 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)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
RNA backbone	8273	3508	-
Q-score	-	25397	14717 (2.90 - 3.90)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the EM map (all-atom inclusion $< 40\%$). The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	B7	95	
2	B8	188	
3	B9	100	

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Mol	Chain	Length	Quality of chain
4	BA	1571	
5	BB	73	
6	BD	306	
7	BE	399	
8	BF	294	
9	BI	268	
10	BJ	262	
11	BK	192	
12	BL	184	
13	BN	178	
14	BO	145	
15	BP	296	
16	BQ	251	
17	BR	169	
18	BS	180	
19	BT	292	
20	BU	149	
21	BV	209	
22	BW	210	
23	BX	150	
24	BY	216	
25	Ba	423	
26	Bb	380	
27	Bc	334	
28	Bd	206	

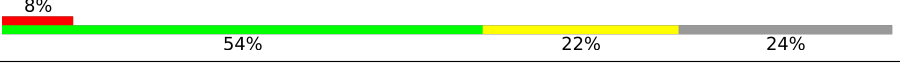
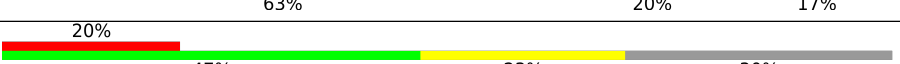
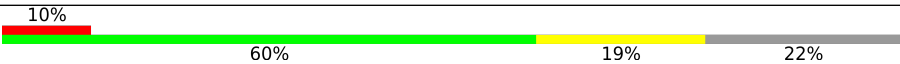
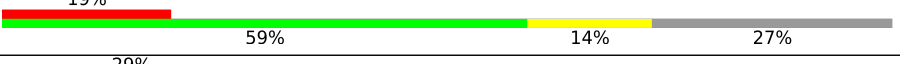
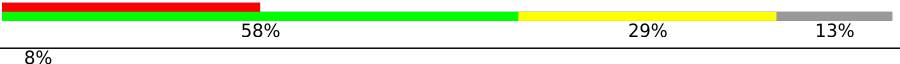
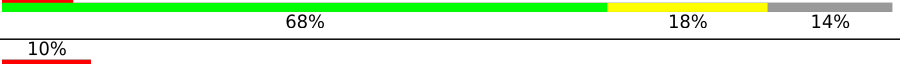
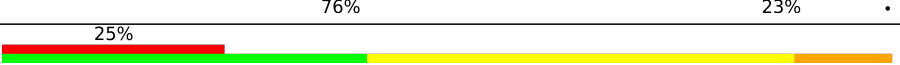
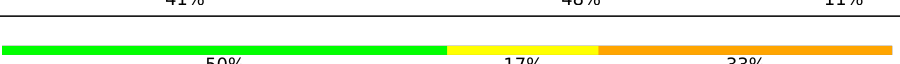
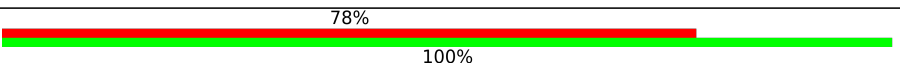
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Mol	Chain	Length	Quality of chain
29	Be	135	
30	Bf	142	
31	Bg	159	
32	Bh	332	
33	Bi	306	
34	Bj	279	
35	Bk	269	
36	Bl	166	
37	Bm	198	
38	Bn	128	
39	Bo	124	
40	Bp	112	
41	Bq	138	
42	Bt	102	
43	Bu	205	
44	Bv	222	
45	Bw	433	
46	Bx	196	
47	AA	962	
48	CL	198	
48	DL	198	
48	EL	198	
48	FL	198	
48	GL	198	
48	HL	198	

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Mol	Chain	Length	Quality of chain
49	AB	289	
50	AC	167	
51	AE	430	
52	AF	276	
53	AG	242	
54	AI	397	
55	AJ	200	
56	AK	196	
57	AL	139	
58	AN	128	
59	AO	239	
60	AP	135	
61	AQ	130	
62	AR	143	
63	AU	87	
64	AV	71	
65	AX	6	
66	AZ	18	
67	Aa	382	
68	Ab	190	
69	Ac	173	
70	Ad	205	
71	Ae	455	
72	Af	188	
73	Ag	410	

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Mol	Chain	Length	Quality of chain
74	Ah	387	
75	Ai	106	
76	Aj	218	
77	Ak	325	
78	Am	118	
79	An	199	
80	Ao	699	
81	Ap	258	
82	B0	148	
83	B1	256	
84	B2	252	
85	B3	161	
86	B4	126	
87	B5	188	
88	B6	65	

2 Entry composition [i](#)

There are 94 unique types of molecules in this entry. The entry contains 174319 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 protein called Mitochondrial ribosomal protein L34.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	B7	46	Total	C	N	O	S	0	0
			387	239	89	58	1		

- Molecule 2 is a protein called Mitochondrial ribosomal protein L35.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B8	95	Total	C	N	O	S	0	0
			833	539	163	129	2		

- Molecule 3 is a protein called Ribosomal protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	B9	38	Total	C	N	O	S	0	0
			335	214	70	47	4		

- Molecule 4 is a RNA chain called 16S rRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	BA	1544	Total	C	N	O	P	0	0
			32844	14750	5972	10578	1544		

- Molecule 5 is a RNA chain called CP tRNAPhe.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	BB	67	Total	C	N	O	P	0	0
			1427	640	261	459	67		

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
BB	72	C	-	insertion	GB 76262549
BB	73	A	-	insertion	GB 76262549

- Molecule 6 is a protein called uL2m.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	BD	240	Total	C	N	O	S	0	0
			1860	1160	371	319	10		

- Molecule 7 is a protein called ICT1.

Mol	Chain	Residues	Atoms					AltConf	Trace
7	BE	307	Total	C	N	O	S	0	0
			2420	1554	426	430	10		

- Molecule 8 is a protein called Mitochondrial ribosomal protein L4.

Mol	Chain	Residues	Atoms					AltConf	Trace
8	BF	250	Total	C	N	O	S	0	0
			2011	1294	367	344	6		

- Molecule 9 is a protein called Mitochondrial ribosomal protein L9.

Mol	Chain	Residues	Atoms				AltConf	Trace
9	BI	98	Total	C	N	O	0	0
			805	509	155	141		

- Molecule 10 is a protein called Mitochondrial ribosomal protein L10.

Mol	Chain	Residues	Atoms					AltConf	Trace
10	BJ	212	Total	C	N	O	S	0	0
			1705	1100	306	290	9		

- Molecule 11 is a protein called Mitochondrial ribosomal protein L11.

Mol	Chain	Residues	Atoms					AltConf	Trace
11	BK	176	Total	C	N	O	S	0	0
			1303	830	236	235	2		

- Molecule 12 is a protein called Peptidyl-tRNA hydrolase ICT1, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
12	BL	137	Total	C	N	O	S	0	0
			1113	683	219	206	5		

There are 7 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
BL	23	SER	-	expression tag	UNP Q14197
BL	24	GLY	-	expression tag	UNP Q14197
BL	25	GLY	-	expression tag	UNP Q14197
BL	26	SER	-	expression tag	UNP Q14197
BL	27	GLY	-	expression tag	UNP Q14197
BL	28	SER	-	expression tag	UNP Q14197
BL	29	GLY	-	expression tag	UNP Q14197

- Molecule 13 is a protein called uL13m.

Mol	Chain	Residues	Atoms					AltConf	Trace
13	BN	177	Total	C	N	O	S	0	0
			1444	926	258	253	7		

- Molecule 14 is a protein called uL14m.

Mol	Chain	Residues	Atoms					AltConf	Trace
14	BO	115	Total	C	N	O	S	0	0
			896	562	176	154	4		

- Molecule 15 is a protein called uL15m.

Mol	Chain	Residues	Atoms					AltConf	Trace
15	BP	288	Total	C	N	O	S	0	0
			2312	1473	430	403	6		

- Molecule 16 is a protein called uL16m.

Mol	Chain	Residues	Atoms					AltConf	Trace
16	BQ	222	Total	C	N	O	S	0	0
			1803	1156	331	306	10		

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
BQ	237	HIS	TYR	conflict	UNP F1RI89

- Molecule 17 is a protein called bL17m.

Mol	Chain	Residues	Atoms					AltConf	Trace
17	BR	153	Total	C	N	O	S	0	0
			1240	777	236	222	5		

- Molecule 18 is a protein called Mitochondrial ribosomal protein L18.

Mol	Chain	Residues	Atoms					AltConf	Trace
18	BS	143	Total	C	N	O	S	0	0
			1168	733	227	204	4		

- Molecule 19 is a protein called Mitochondrial ribosomal protein L19.

Mol	Chain	Residues	Atoms					AltConf	Trace
19	BT	240	Total	C	N	O	S	0	0
			1954	1253	338	354	9		

- Molecule 20 is a protein called Mitochondrial ribosomal protein L20.

Mol	Chain	Residues	Atoms					AltConf	Trace
20	BU	140	Total	C	N	O	S	0	0
			1159	732	239	185	3		

- Molecule 21 is a protein called Mitochondrial ribosomal protein L21.

Mol	Chain	Residues	Atoms					AltConf	Trace
21	BV	155	Total	C	N	O	S	0	0
			1231	789	219	219	4		

- Molecule 22 is a protein called uL22m.

Mol	Chain	Residues	Atoms					AltConf	Trace
22	BW	166	Total	C	N	O	S	0	0
			1374	876	258	234	6		

- Molecule 23 is a protein called uL23m.

Mol	Chain	Residues	Atoms					AltConf	Trace
23	BX	149	Total	C	N	O	S	0	0
			1181	752	227	200	2		

- Molecule 24 is a protein called uL24m.

Mol	Chain	Residues	Atoms					AltConf	Trace
24	BY	206	Total	C	N	O	S	0	0
			1678	1056	308	309	5		

- Molecule 25 is a protein called Mitochondrial ribosomal protein L37.

Mol	Chain	Residues	Atoms					AltConf	Trace
25	Ba	393	Total	C	N	O	S	0	0
			3173	2040	556	565	12		

- Molecule 26 is a protein called Mitochondrial ribosomal protein L38.

Mol	Chain	Residues	Atoms					AltConf	Trace
26	Bb	354	Total	C	N	O	S	0	0
			2952	1876	542	525	9		

- Molecule 27 is a protein called Mitochondrial ribosomal protein L39.

Mol	Chain	Residues	Atoms					AltConf	Trace
27	Bc	295	Total	C	N	O	S	0	0
			2408	1541	410	441	16		

- Molecule 28 is a protein called Mitochondrial ribosomal protein L40.

Mol	Chain	Residues	Atoms					AltConf	Trace
28	Bd	138	Total	C	N	O	S	0	0
			1158	729	211	217	1		

- Molecule 29 is a protein called Mitochondrial ribosomal protein L41.

Mol	Chain	Residues	Atoms					AltConf	Trace
29	Be	122	Total	C	N	O	S	0	0
			972	628	168	173	3		

- Molecule 30 is a protein called mL42.

Mol	Chain	Residues	Atoms					AltConf	Trace
30	Bf	108	Total	C	N	O	S	0	0
			827	519	154	150	4		

- Molecule 31 is a protein called Mitochondrial ribosomal protein L43.

Mol	Chain	Residues	Atoms					AltConf	Trace
31	Bg	148	Total	C	N	O	S	0	0
			1167	727	225	212	3		

- Molecule 32 is a protein called mL44.

Mol	Chain	Residues	Atoms					AltConf	Trace
32	Bh	289	Total	C	N	O	S	0	0
			2319	1486	399	426	8		

- Molecule 33 is a protein called Mitochondrial ribosomal protein L45.

Mol	Chain	Residues	Atoms					AltConf	Trace
33	Bi	260	Total	C	N	O	S	0	0
			2138	1370	379	379	10		

- Molecule 34 is a protein called Mitochondrial ribosomal protein L46.

Mol	Chain	Residues	Atoms					AltConf	Trace
34	Bj	217	Total	C	N	O	S	0	0
			1775	1137	311	321	6		

- Molecule 35 is a protein called 39S ribosomal protein L48, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
35	Bk	155	Total	C	N	O	S	0	0
			1246	796	214	231	5		

- Molecule 36 is a protein called Mrpl34.

Mol	Chain	Residues	Atoms					AltConf	Trace
36	Bl	133	Total	C	N	O	S	0	0
			1097	709	192	194	2		

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
Bl	59	ARG	LYS	conflict	UNP A0A0R4J8D6

- Molecule 37 is a protein called Mitochondrial ribosomal protein L50.

Mol	Chain	Residues	Atoms					AltConf	Trace
37	Bm	109	Total	C	N	O	S	0	0
			893	568	160	162	3		

- Molecule 38 is a protein called Mitochondrial ribosomal protein L51.

Mol	Chain	Residues	Atoms					AltConf	Trace
38	Bn	97	Total	C	N	O	S	0	0
			837	539	166	128	4		

- Molecule 39 is a protein called mL52.

Mol	Chain	Residues	Atoms					AltConf	Trace
39	Bo	97	Total	C	N	O	S	0	0
			772	481	148	141	2		

- Molecule 40 is a protein called mL53.

Mol	Chain	Residues	Atoms					AltConf	Trace
40	Bp	97	Total	C	N	O	S	0	0
			742	459	143	134	6		

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
Bp	12	ALA	SER	conflict	UNP A0A341D604
Bp	107	SER	GLY	conflict	UNP A0A341D604

- Molecule 41 is a protein called mL54.

Mol	Chain	Residues	Atoms					AltConf	Trace
41	Bq	54	Total	C	N	O	S	0	0
			474	304	88	81	1		

- Molecule 42 is a protein called Mitochondrial ribosomal protein L57.

Mol	Chain	Residues	Atoms					AltConf	Trace
42	Bt	94	Total	C	N	O	S	0	0
			780	485	168	126	1		

- Molecule 43 is a protein called mL62 (ICT1).

Mol	Chain	Residues	Atoms					AltConf	Trace
43	Bu	151	Total	C	N	O	S	0	0
			1198	738	233	222	5		

- Molecule 44 is a protein called mL64.

Mol	Chain	Residues	Atoms					AltConf	Trace
44	Bv	135	Total	C	N	O	S	0	0
			1131	692	223	211	5		

- Molecule 45 is a protein called 39S ribosomal protein S30, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
45	Bw	387	Total	C	N	O	S	0	0
			3126	2011	548	555	12		

- Molecule 46 is a protein called Mitochondrial ribosomal protein S18A.

Mol	Chain	Residues	Atoms					AltConf	Trace
46	Bx	162	Total	C	N	O	S	0	0
			1325	845	249	224	7		

- Molecule 47 is a RNA chain called 12S rRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
47	AA	960	Total	C	N	O	P	0	0
			20411	9162	3708	6581	960		

- Molecule 48 is a protein called Mitochondrial ribosomal protein L12.

Mol	Chain	Residues	Atoms				AltConf	Trace
48	CL	45	Total	C	N	O	0	0
			317	203	52	62		
48	DL	27	Total	C	N	O	0	0
			213	137	33	43		
48	EL	28	Total	C	N	O	0	0
			222	143	35	44		
48	FL	27	Total	C	N	O	0	0
			213	137	33	43		
48	GL	27	Total	C	N	O	0	0
			213	137	33	43		
48	HL	26	Total	C	N	O	0	0
			205	131	32	42		

- Molecule 49 is a protein called Mitochondrial ribosomal protein S2.

Mol	Chain	Residues	Atoms					AltConf	Trace
49	AB	220	Total	C	N	O	S	0	0
			1762	1126	326	304	6		

- Molecule 50 is a protein called Mitochondrial ribosomal protein S24.

Mol	Chain	Residues	Atoms					AltConf	Trace
50	AC	132	Total	C	N	O	S	0	0
			1075	695	195	181	4		

- Molecule 51 is a protein called Mitochondrial ribosomal protein S5.

Mol	Chain	Residues	Atoms					AltConf	Trace
51	AE	343	Total	C	N	O	S	0	0
			2732	1707	527	487	11		

- Molecule 52 is a protein called bS6m.

Mol	Chain	Residues	Atoms					AltConf	Trace
52	AF	122	Total	C	N	O	S	0	0
			981	620	178	177	6		

- Molecule 53 is a protein called Mitochondrial ribosomal protein S7.

Mol	Chain	Residues	Atoms					AltConf	Trace
53	AG	208	Total	C	N	O	S	0	0
			1721	1097	314	299	11		

- Molecule 54 is a protein called uS9m.

Mol	Chain	Residues	Atoms					AltConf	Trace
54	AI	328	Total	C	N	O	S	0	0
			2650	1678	478	481	13		

- Molecule 55 is a protein called Mitochondrial ribosomal protein S10.

Mol	Chain	Residues	Atoms					AltConf	Trace
55	AJ	140	Total	C	N	O	S	0	0
			1155	746	197	208	4		

- Molecule 56 is a protein called uS11m.

Mol	Chain	Residues	Atoms					AltConf	Trace
56	AK	137	Total	C	N	O	S	0	0
			1007	631	193	180	3		

- Molecule 57 is a protein called Mitochondrial ribosomal protein S12.

Mol	Chain	Residues	Atoms					AltConf	Trace
57	AL	109	Total	C	N	O	S	0	0
			840	524	172	138	6		

- Molecule 58 is a protein called Mitochondrial ribosomal protein S14.

Mol	Chain	Residues	Atoms					AltConf	Trace
58	AN	101	Total	C	N	O	S	0	0
			858	534	174	144	6		

- Molecule 59 is a protein called uS15m.

Mol	Chain	Residues	Atoms					AltConf	Trace
59	AO	175	Total	C	N	O	S	0	0
			1448	919	272	248	9		

- Molecule 60 is a protein called 28S ribosomal protein S16, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
60	AP	117	Total	C	N	O	S	0	0
			932	588	184	155	5		

- Molecule 61 is a protein called uS17m.

Mol	Chain	Residues	Atoms					AltConf	Trace
61	AQ	112	Total	C	N	O	S	0	0
			875	568	153	151	3		

- Molecule 62 is a protein called Mitochondrial ribosomal protein S18C.

Mol	Chain	Residues	Atoms					AltConf	Trace
62	AR	97	Total	C	N	O	S	0	0
			784	507	132	138	7		

- Molecule 63 is a protein called bS21m.

Mol	Chain	Residues	Atoms					AltConf	Trace
63	AU	86	Total	C	N	O	S	0	0
			734	453	148	125	8		

- Molecule 64 is a RNA chain called fMet-tRNAMet (P site).

Mol	Chain	Residues	Atoms					AltConf	Trace
64	AV	71	Total	C	N	O	P	0	0
			1498	673	264	491	70		

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
AV	69	C	-	insertion	GB 1208989970
AV	70	C	-	insertion	GB 1208989970
AV	71	A	-	insertion	GB 1208989970

- Molecule 65 is a RNA chain called mRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
65	AX	6	Total	C	N	O	P	0	0
			131	59	27	39	6		

- Molecule 66 is a protein called unknown.

Mol	Chain	Residues	Atoms				AltConf	Trace
66	AZ	18	Total	C	N	O	0	0
			90	54	18	18		

- Molecule 67 is a protein called Mitochondrial ribosomal protein S22.

Mol	Chain	Residues	Atoms					AltConf	Trace
67	Aa	292	Total	C	N	O	S	0	0
			2378	1518	409	442	9		

- Molecule 68 is a protein called mS23.

Mol	Chain	Residues	Atoms					AltConf	Trace
68	Ab	135	Total	C	N	O	S	0	0
			1101	709	199	192	1		

- Molecule 69 is a protein called Mitochondrial ribosomal protein S25.

Mol	Chain	Residues	Atoms					AltConf	Trace
69	Ac	169	Total	C	N	O	S	0	0
			1367	876	236	245	10		

- Molecule 70 is a protein called Mitochondrial ribosomal protein S26.

Mol	Chain	Residues	Atoms					AltConf	Trace
70	Ad	177	Total	C	N	O	S	0	0
			1467	904	288	273	2		

- Molecule 71 is a protein called Mitochondrial ribosomal protein S27.

Mol	Chain	Residues	Atoms					AltConf	Trace
71	Ae	388	Total	C	N	O	S	0	0
			3109	1971	535	589	14		

- Molecule 72 is a protein called Mitoribosomal protein ms28, mrps28.

Mol	Chain	Residues	Atoms					AltConf	Trace
72	Af	99	Total	C	N	O	S	0	0
			778	494	134	146	4		

- Molecule 73 is a protein called Death associated protein 3.

Mol	Chain	Residues	Atoms					AltConf	Trace
73	Ag	353	Total	C	N	O	S	0	0
			2875	1837	515	513	10		

- Molecule 74 is a protein called Mitochondrial ribosomal protein S31.

Mol	Chain	Residues	Atoms					AltConf	Trace
74	Ah	120	Total	C	N	O	S	0	0
			1015	659	168	185	3		

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
Ah	180	UNK	-	insertion	UNP F1RME2
Ah	181	GLN	-	insertion	UNP F1RME2
Ah	182	LYS	-	insertion	UNP F1RME2
Ah	184	GLY	-	insertion	UNP F1RME2
Ah	185	GLU	LYS	conflict	UNP F1RME2
Ah	187	PRO	LYS	conflict	UNP F1RME2
Ah	189	ILE	LEU	conflict	UNP F1RME2
Ah	190	SER	ILE	conflict	UNP F1RME2
Ah	237	SER	-	insertion	UNP F1RME2
Ah	238	PHE	-	insertion	UNP F1RME2

- Molecule 75 is a protein called mS33.

Mol	Chain	Residues	Atoms					AltConf	Trace
75	Ai	99	Total	C	N	O	S	0	0
			824	522	156	143	3		

- Molecule 76 is a protein called mS34.

Mol	Chain	Residues	Atoms					AltConf	Trace
76	Aj	213	Total	C	N	O	S	0	0
			1788	1131	338	311	8		

- Molecule 77 is a protein called Mitochondrial ribosomal protein S35.

Mol	Chain	Residues	Atoms					AltConf	Trace
77	Ak	275	Total	C	N	O	S	0	0
			2222	1414	380	419	9		

- Molecule 78 is a protein called mS37.

Mol	Chain	Residues	Atoms					AltConf	Trace
78	Am	116	Total	C	N	O	S	0	0
			930	577	185	160	8		

- Molecule 79 is a protein called Aurora kinase A interacting protein 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
79	An	72	Total	C	N	O	S	0	0
			639	407	139	92	1		

- Molecule 80 is a protein called Pentatricopeptide repeat domain 3.

Mol	Chain	Residues	Atoms					AltConf	Trace
80	Ao	572	Total	C	N	O	S	0	0
			4526	2898	770	834	24		

- Molecule 81 is a protein called 28S ribosomal protein S18b, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
81	Ap	190	Total	C	N	O	S	0	0
			1564	991	292	273	8		

- Molecule 82 is a protein called Mitochondrial ribosomal protein L27.

Mol	Chain	Residues	Atoms					AltConf	Trace
82	B0	110	Total	C	N	O	S	0	0
			857	553	156	145	3		

- Molecule 83 is a protein called Mitochondrial ribosomal protein L28.

Mol	Chain	Residues	Atoms					AltConf	Trace
83	B1	244	Total	C	N	O	S	0	0
			2036	1315	363	353	5		

- Molecule 84 is a protein called Mitochondrial ribosomal protein L47.

Mol	Chain	Residues	Atoms					AltConf	Trace
84	B2	179	Total	C	N	O	S	0	0
			1548	992	290	260	6		

- Molecule 85 is a protein called uL30m.

Mol	Chain	Residues	Atoms					AltConf	Trace
85	B3	118	Total	C	N	O	S	0	0
			968	622	178	165	3		

- Molecule 86 is a protein called bL31m.

Mol	Chain	Residues	Atoms					AltConf	Trace
86	B4	62	Total	C	N	O	S	0	0
			474	296	94	81	3		

- Molecule 87 is a protein called bL32m.

Mol	Chain	Residues	Atoms					AltConf	Trace
87	B5	110	Total	C	N	O	S	0	0
			902	553	181	162	6		

- Molecule 88 is a protein called bL33m.

Mol	Chain	Residues	Atoms					AltConf	Trace
88	B6	52	Total	C	N	O	S	0	0
			425	274	78	71	2		

- Molecule 89 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		AltConf
89	B9	1	Total 1	Zn 1	0
89	Bx	1	Total 1	Zn 1	0
89	AR	1	Total 1	Zn 1	0
89	Ac	1	Total 1	Zn 1	0
89	Ap	1	Total 1	Zn 1	0
89	B5	1	Total 1	Zn 1	0

- Molecule 90 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

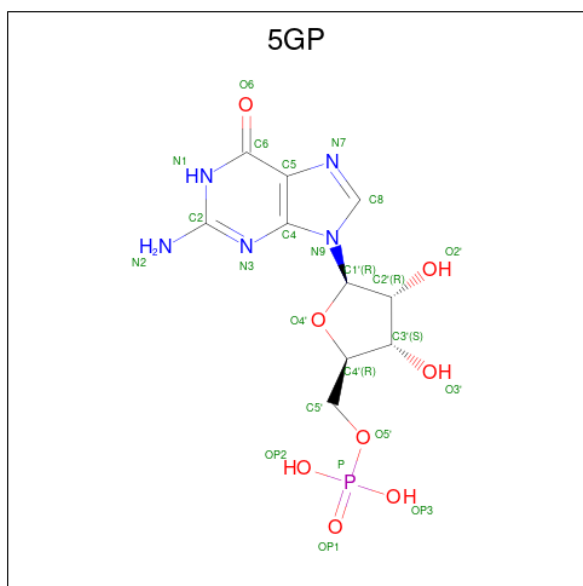
Mol	Chain	Residues	Atoms		AltConf
90	BA	212	Total 212	Mg 212	0
90	BB	1	Total 1	Mg 1	0
90	BD	3	Total 3	Mg 3	0
90	BE	1	Total 1	Mg 1	0
90	BP	3	Total 3	Mg 3	0
90	BQ	1	Total 1	Mg 1	0
90	BR	1	Total 1	Mg 1	0
90	Be	2	Total 2	Mg 2	0
90	Bl	1	Total 1	Mg 1	0
90	Bt	1	Total 1	Mg 1	0
90	AA	101	Total 101	Mg 101	0
90	AB	1	Total 1	Mg 1	0
90	AL	1	Total 1	Mg 1	0
90	AV	1	Total 1	Mg 1	0

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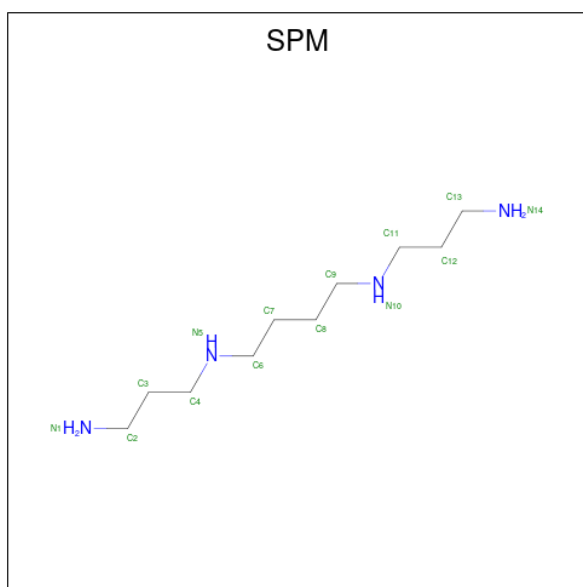
Mol	Chain	Residues	Atoms		AltConf
90	AX	1	Total	Mg	0
			1	1	
90	Ag	1	Total	Mg	0
			1	1	
90	Am	1	Total	Mg	0
			1	1	
90	An	1	Total	Mg	0
			1	1	
90	B3	1	Total	Mg	0
			1	1	

- Molecule 91 is GUANOSINE-5'-MONOPHOSPHATE (CCD ID: 5GP) (formula: $C_{10}H_{14}N_5O_8P$).



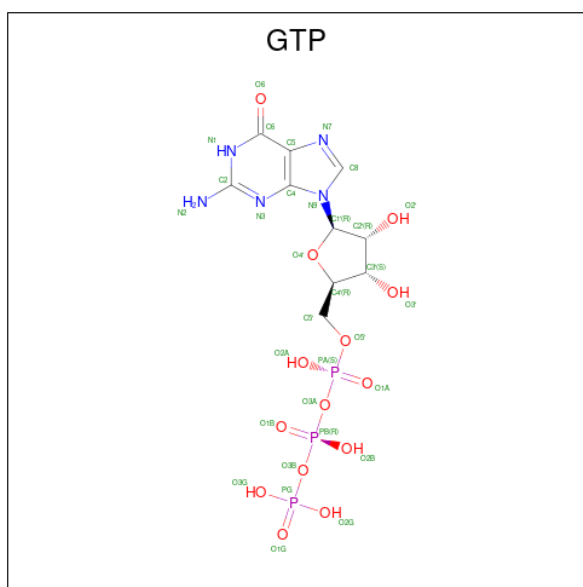
Mol	Chain	Residues	Atoms					AltConf
91	BA	1	Total	C	N	O	P	0
			24	10	5	8	1	
91	BA	1	Total	C	N	O	P	0
			24	10	5	8	1	

- Molecule 92 is SPERMINE (CCD ID: SPM) (formula: $C_{10}H_{26}N_4$).



Mol	Chain	Residues	Atoms			AltConf
92	BA	1	Total 14	C 10	N 4	0
92	BA	1	Total 14	C 10	N 4	0
92	AA	1	Total 14	C 10	N 4	0

- Molecule 93 is GUANOSINE-5'-TRIPHOSPHATE (CCD ID: GTP) (formula: $\text{C}_{10}\text{H}_{16}\text{N}_5\text{O}_{14}\text{P}_3$).

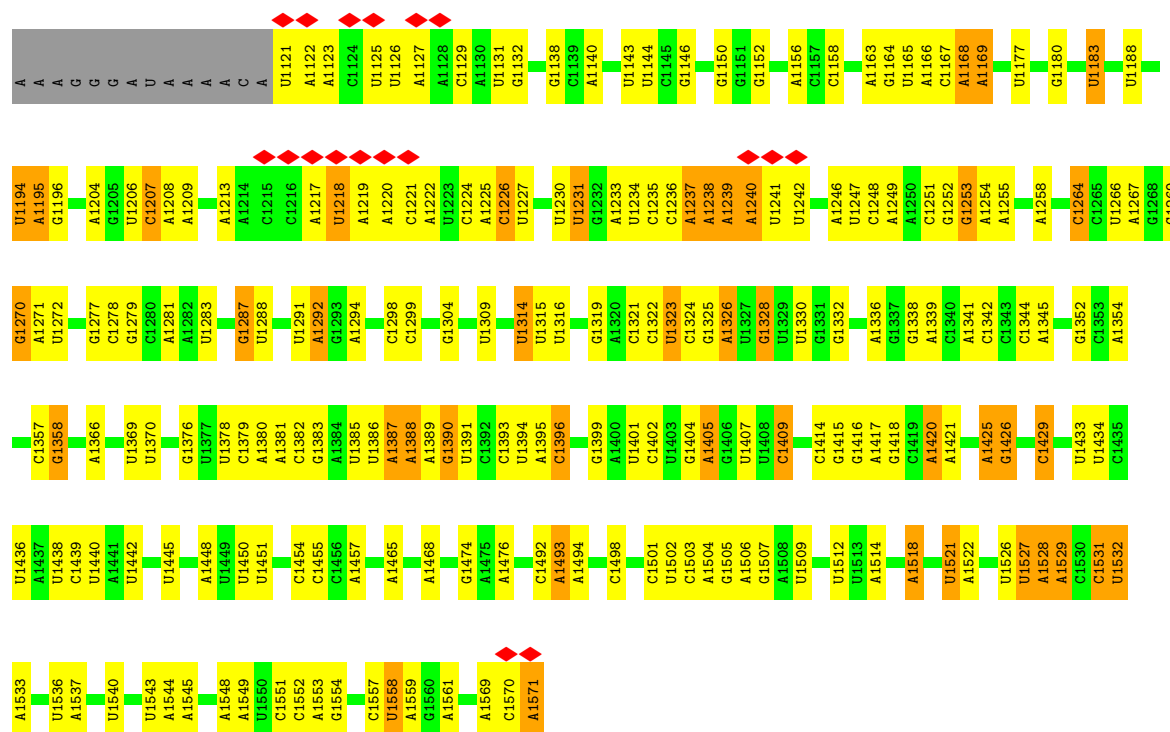


Mol	Chain	Residues	Atoms					AltConf
93	Ag	1	Total	C	N	O	P	0
			32	10	5	14	3	

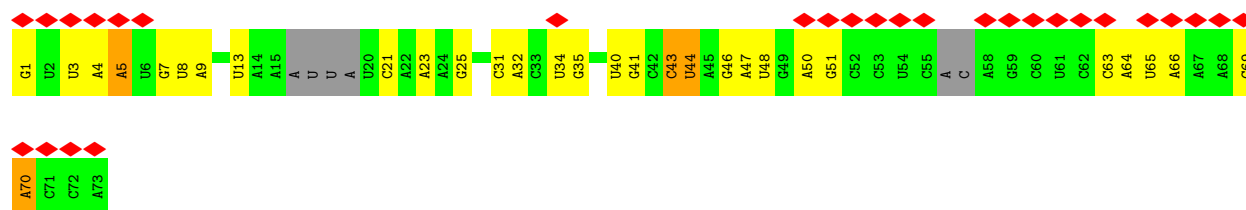
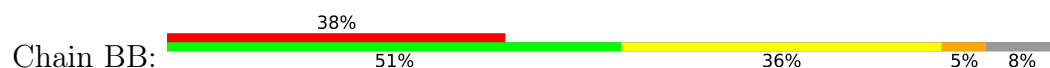
- Molecule 94 is water.

Mol	Chain	Residues	Atoms		AltConf
94	Ag	3	Total	O	0
			3	3	

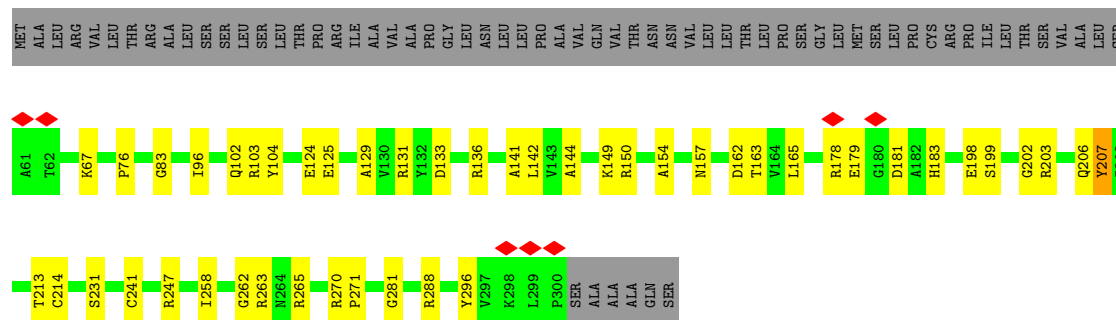




• Molecule 5: CP tRNA^{Phe}



• Molecule 6: uL2m



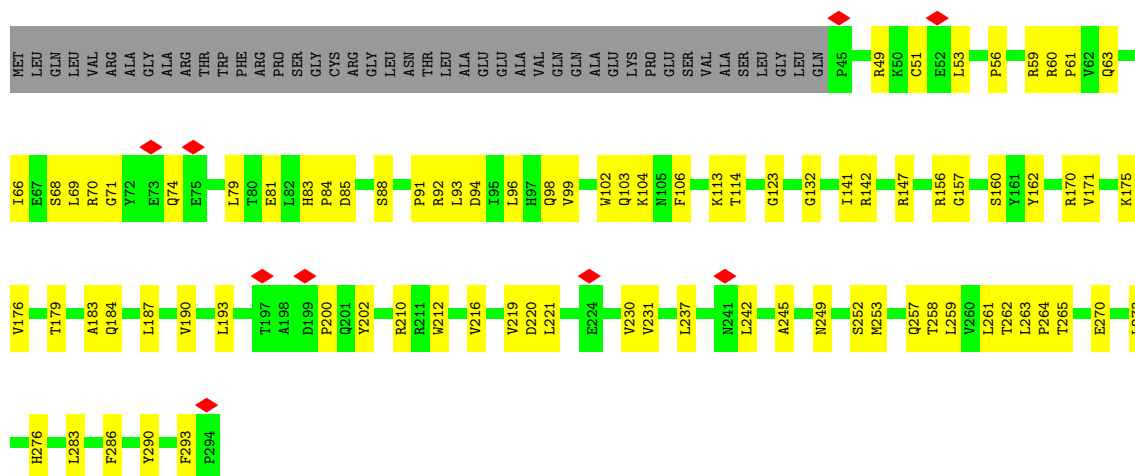
• Molecule 7: ICT1





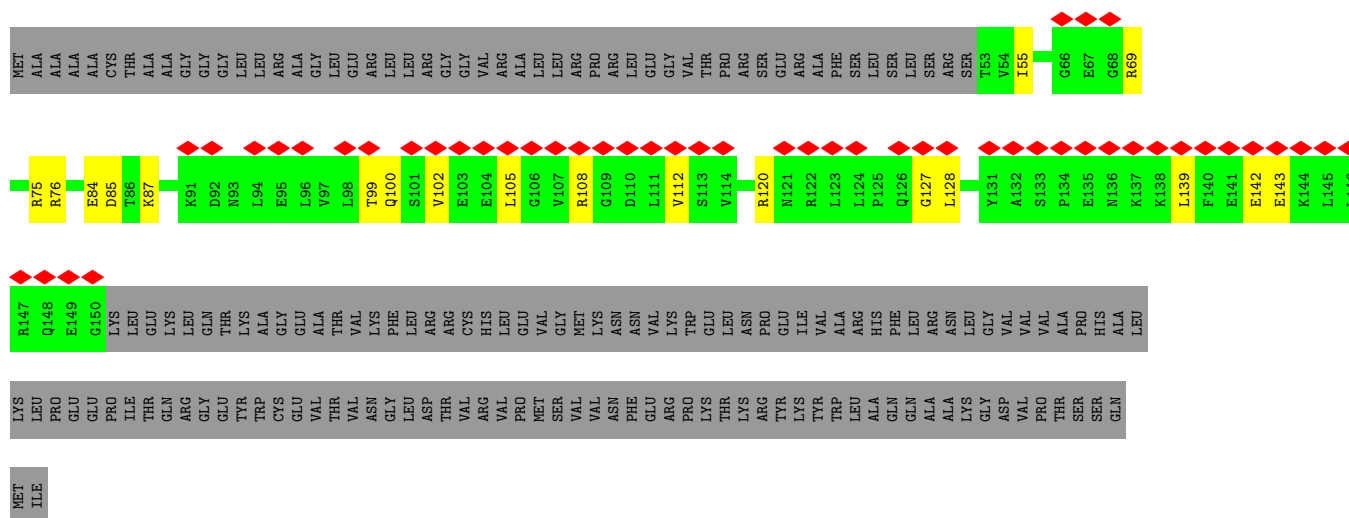
• Molecule 8: Mitochondrial ribosomal protein L4

Chain BF: 57% 28% 15%



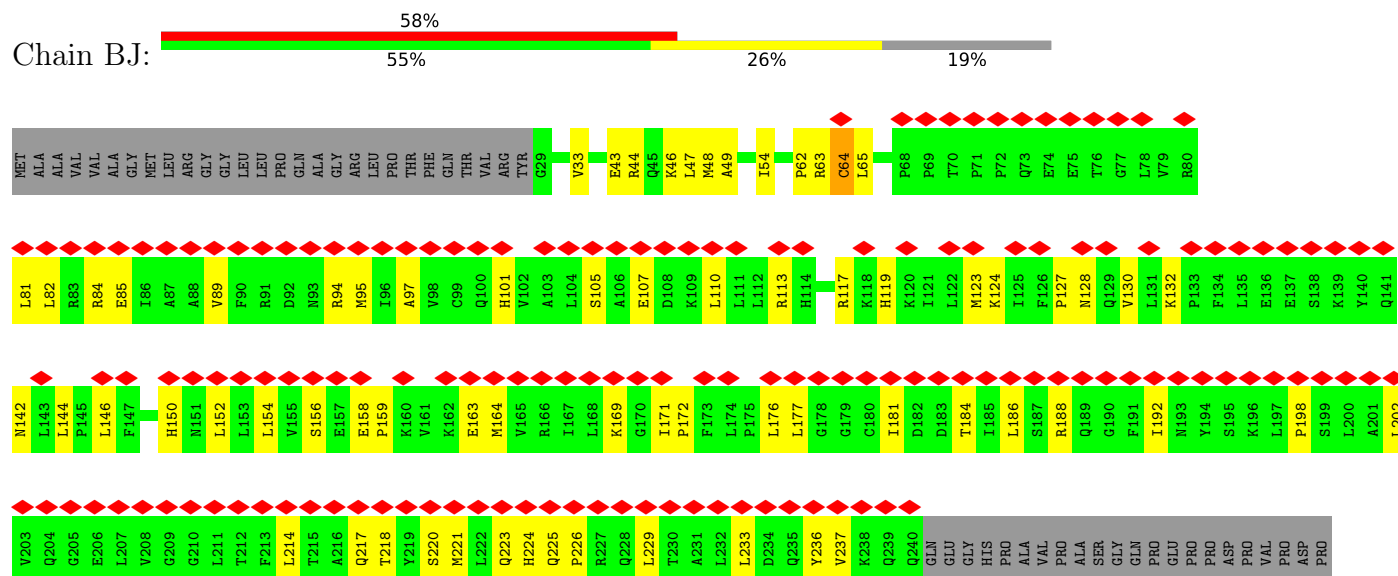
• Molecule 9: Mitochondrial ribosomal protein L9

Chain BI: 19% 29% 7% 63%



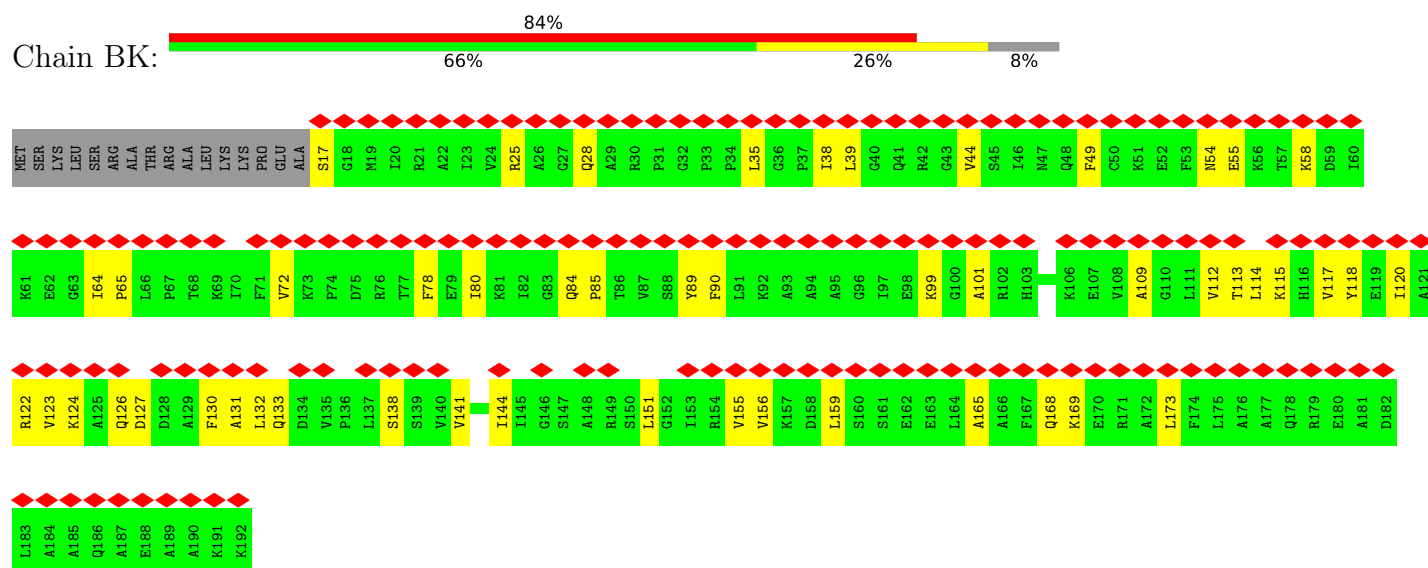
- Molecule 10: Mitochondrial ribosomal protein L10

Chain BJ:



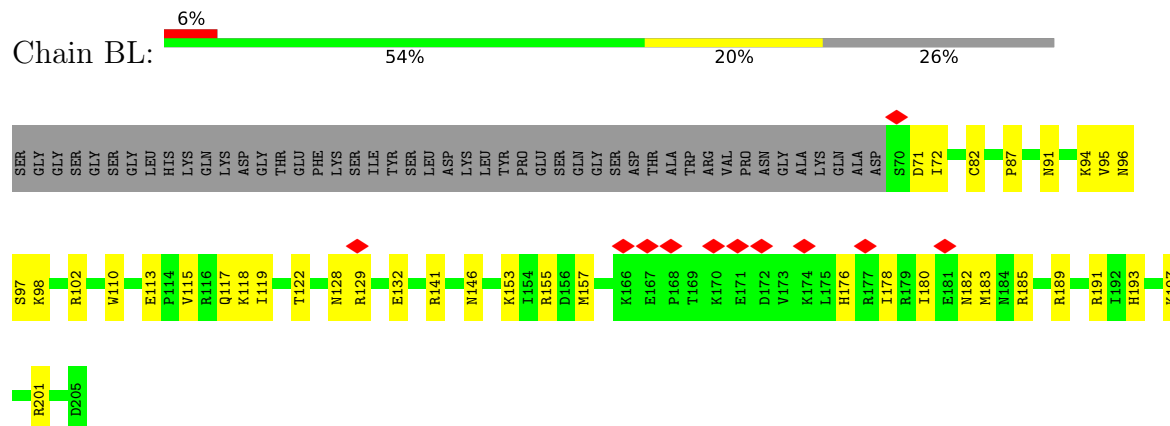
- Molecule 11: Mitochondrial ribosomal protein L11

Chain BK:



- Molecule 12: Peptidyl-tRNA hydrolase ICT1, mitochondrial

Chain BL:



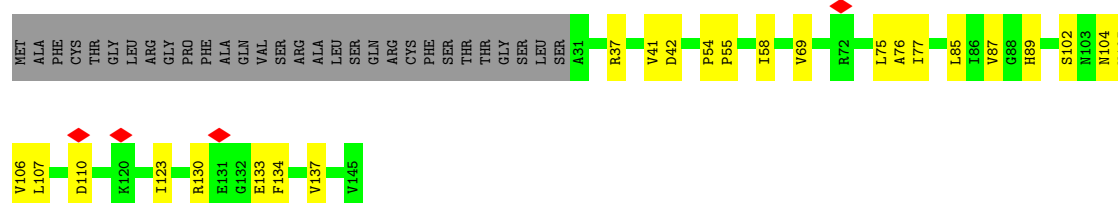
- Molecule 13: uL13m

Chain BN:  75% 24%



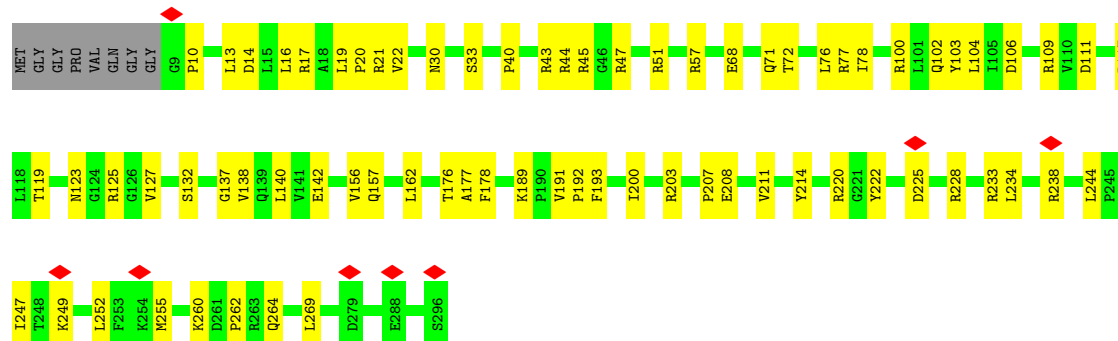
- Molecule 14: uL14m

Chain BO:  63% 17% 21%



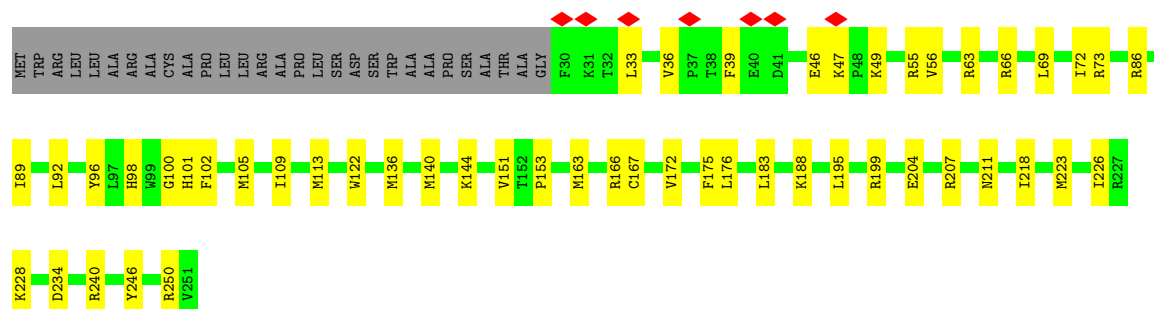
- Molecule 15: uL15m

Chain BP:  72% 25%



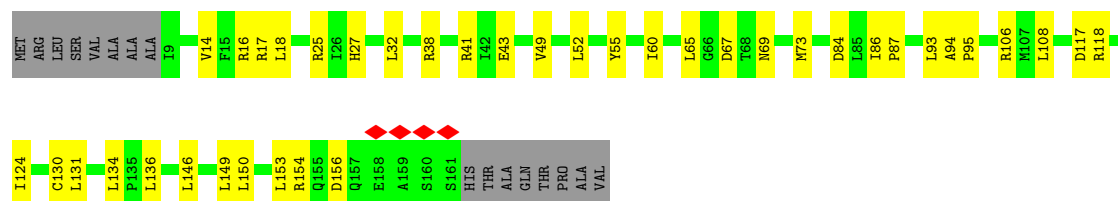
- Molecule 16: uL16m

Chain BQ:  68% 20% 12%



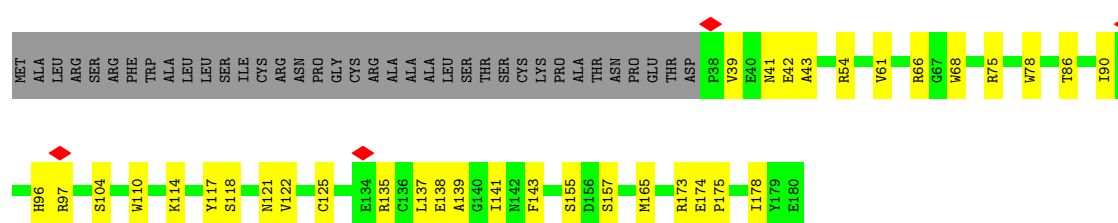
- Molecule 17: bL17m

Chain BR: 



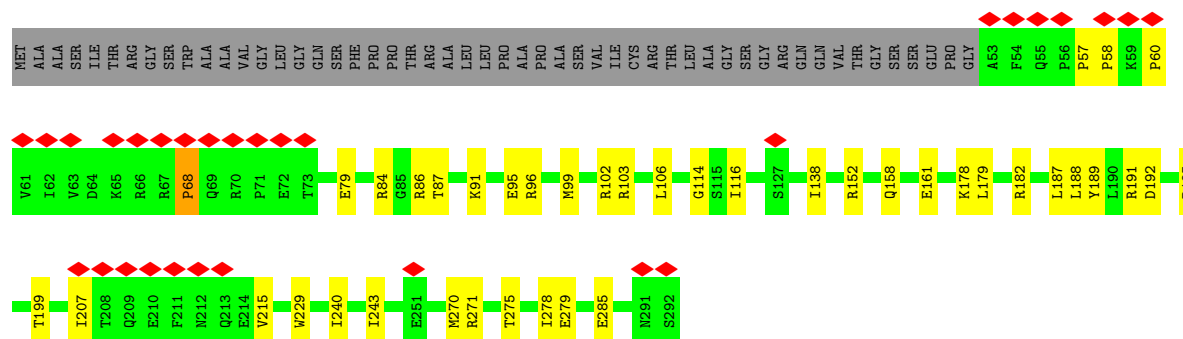
- Molecule 18: Mitochondrial ribosomal protein L18

Chain BS: 



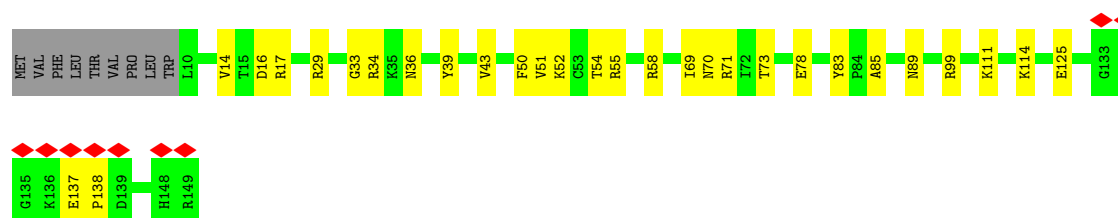
- Molecule 19: Mitochondrial ribosomal protein L19

Chain BT: 



- Molecule 20: Mitochondrial ribosomal protein L20

Chain BU: 



- Molecule 21: Mitochondrial ribosomal protein L21

Sequence logo for the 1000bp upstream region of the human C209 gene. The y-axis represents information content in bits, ranging from 0 to 1.5. The x-axis shows amino acid positions from -1000 to 1999. Amino acids are color-coded: yellow for high frequency, green for medium, and red for low. A red diamond marker is placed above the C209 gene name at position -1000.

The diagram illustrates a network structure with nodes and connections. The nodes are labeled with alphanumeric codes and are color-coded: green, yellow, and grey. The connections are represented by lines between nodes, with some nodes having red diamond markers above them.

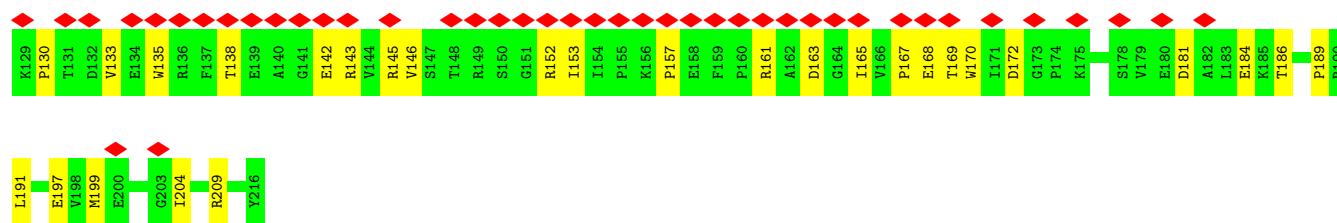
Node Labels and Connections:

- Top Row (Left to Right):** MET (grey), A2 (yellow), V5 (yellow), L6 (green), Y7 (yellow), Y10 (yellow), N14 (yellow), Q27 (yellow), R30 (yellow), E37 (green), T39 (green), V40 (green), Q41 (green), F42 (yellow), R50 (yellow), V51 (yellow), L57 (yellow), A67 (yellow), R71 (green), V72 (green), Q73 (green), H82 (green), R83 (green), R86 (yellow), Q98 (yellow), T104 (yellow), R114 (green), K115 (green), G116 (green), A117 (green), S118 (green), V119 (green), D120 (green), V121 (green), D122 (green), V123 (green), R124 (green), D125 (green), Q126 (green), L128 (green), E129 (green), Q130 (yellow), Q131 (yellow).
- Bottom Row (Left to Right):** R132 (green), Q133 (green), K134 (green), H135 (green), S136 (green), P137 (green), D138 (green), R140 (yellow), R141 (yellow), G142 (green), G143 (green), V144 (green), F145 (green), G146 (green), W147 (green), F148 (green), G149 (green), L150 (yellow).

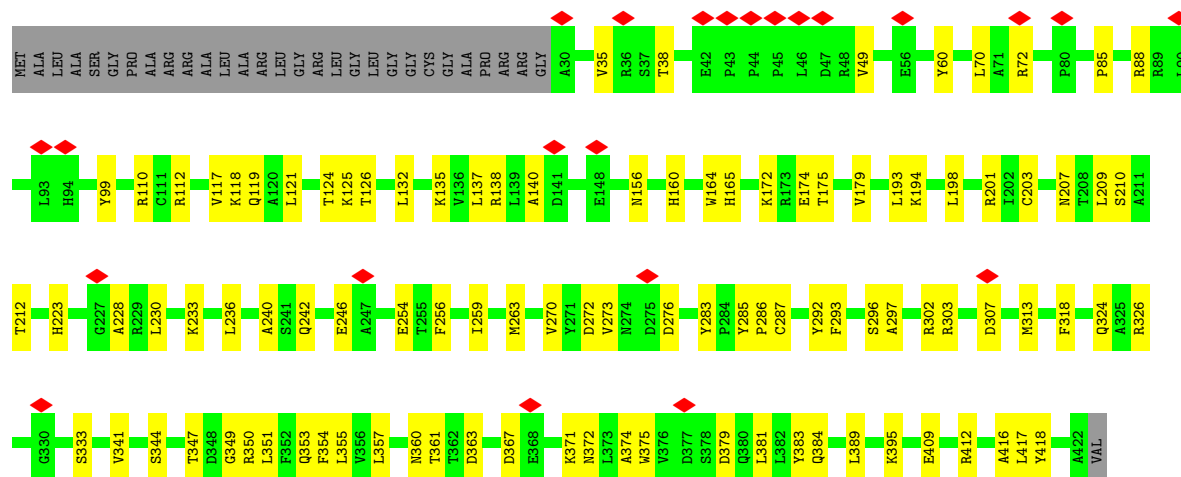
Connections and Markers:

- Red diamond markers are placed above the following nodes: E37, H82, R83, R114, K115, G116, A117, S118, V119, D120, V121, D122, V123, R124, D125, Q126, L128, E129, Q130, Q131.
- Connections are shown as lines between nodes, primarily between adjacent nodes in the same row and between nodes in different rows.

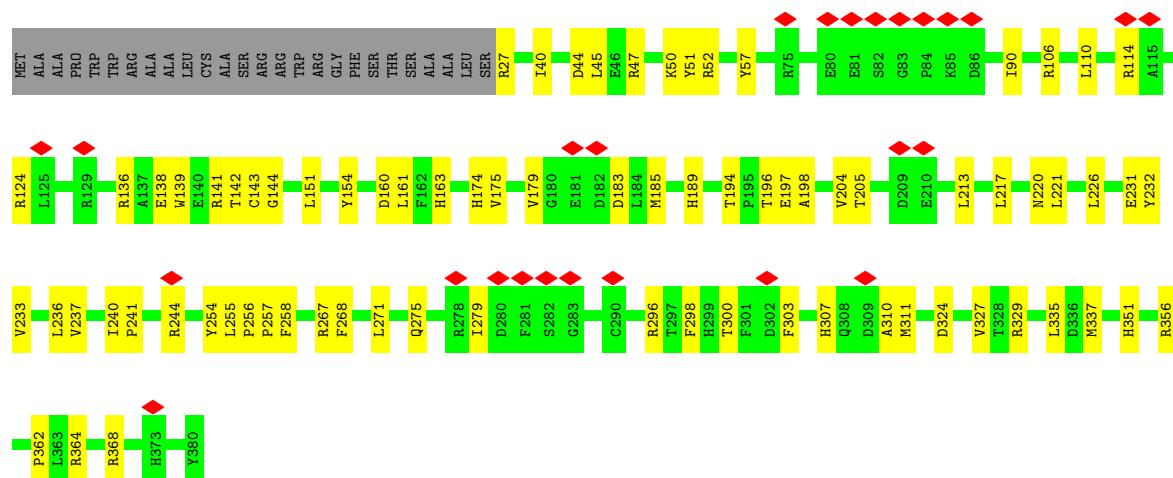
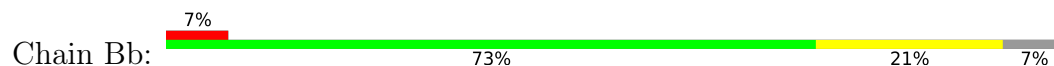
Met	Arg	Leu	Ser	Ala	Leu	Leu	Ala	Leu	Ala	S11	K12	V13	T14	L15	P16	P17	N18	Y19	G22	M23	S24	R25	P26	G27	S28	L29	S30	D31	K32	R33	K34	N35	P36	P37	G38	T39	R40	R43	V44	P45	V46	E47	P48	T49	S50	D51	E52	D53	H54	H55	Q119	V120	K121	L122	V123	P125	T126	D127	L128
E63	I64	L65	E66	G67	K68	D69	A70	G71	K72	V76	V77	Q78	V79	L80	R81	Q82	R83	M84	E89	G90	L91	N92	T93	H94	F95	R96	Y97	V98	G99	K100	T101	V102	D103	Y104	R105	G106	T107	P110	S111	E112	A113	P114	L115	L116	H117	M118	Q119	V120 <td>K121</td> <td>L122</td> <td>V123</td> <td>P125</td> <td>T126</td> <td>D127</td> <td>L128</td>	K121	L122	V123	P125	T126	D127	L128				



• Molecule 25: Mitochondrial ribosomal protein L37

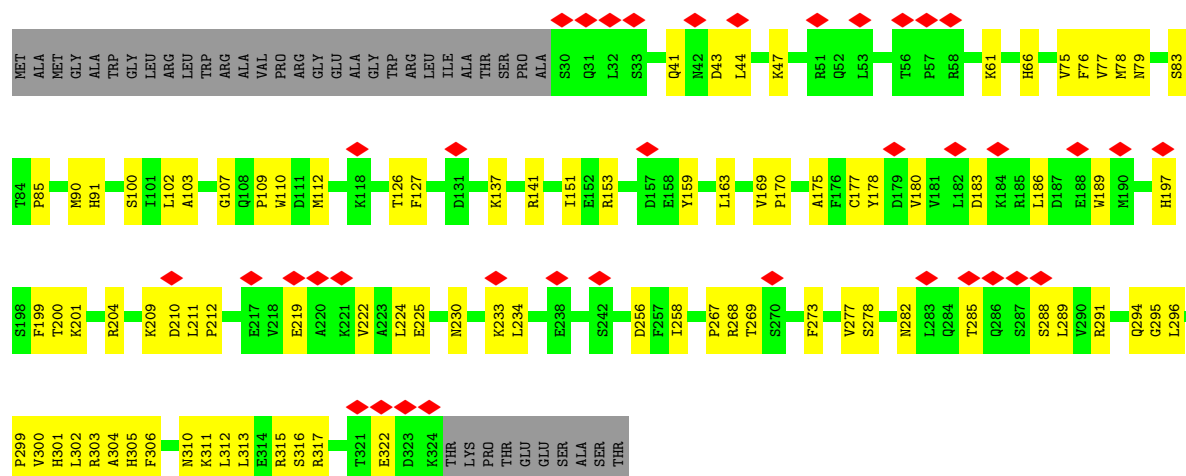


• Molecule 26: Mitochondrial ribosomal protein L38

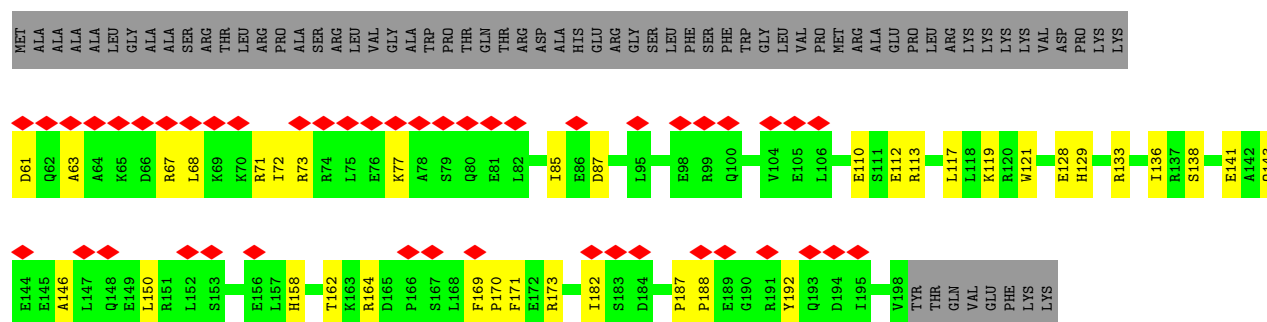


• Molecule 27: Mitochondrial ribosomal protein L39

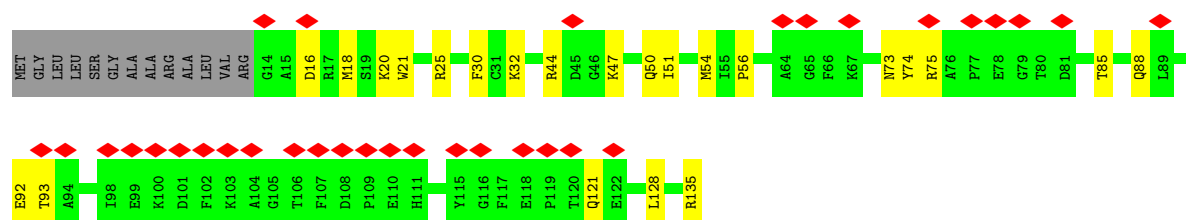
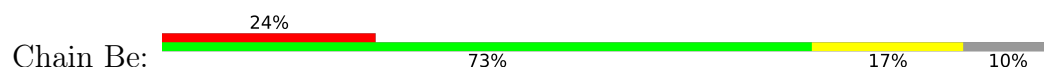




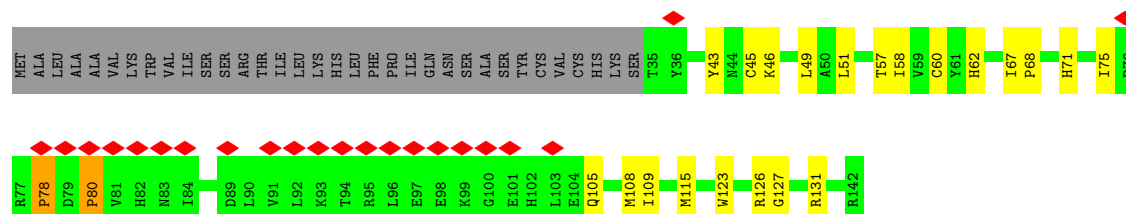
• Molecule 28: Mitochondrial ribosomal protein L40



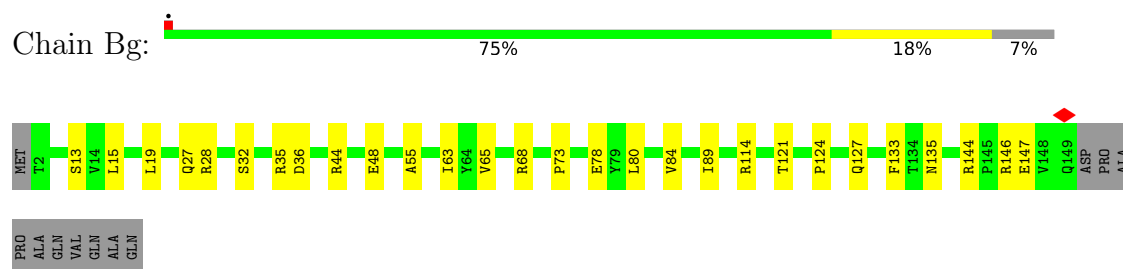
• Molecule 29: Mitochondrial ribosomal protein L41



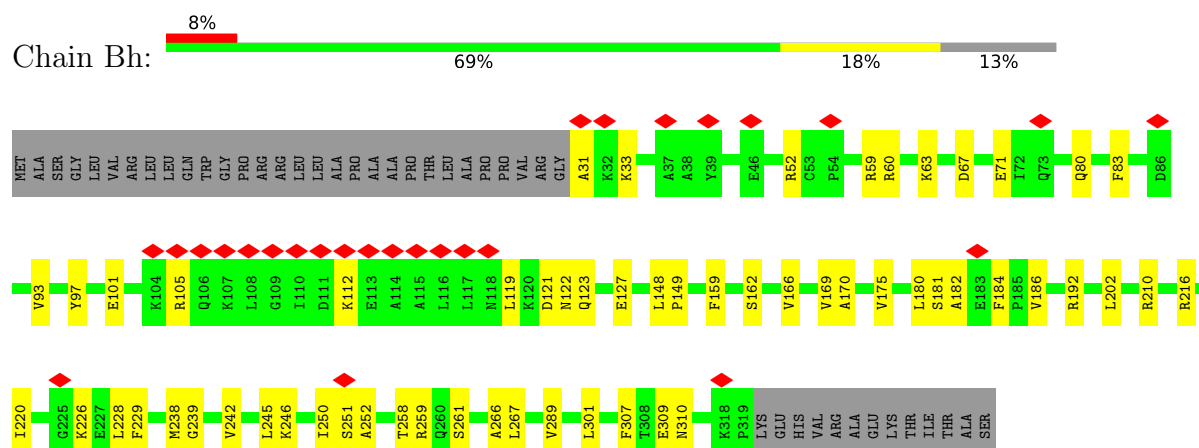
• Molecule 30: mL42



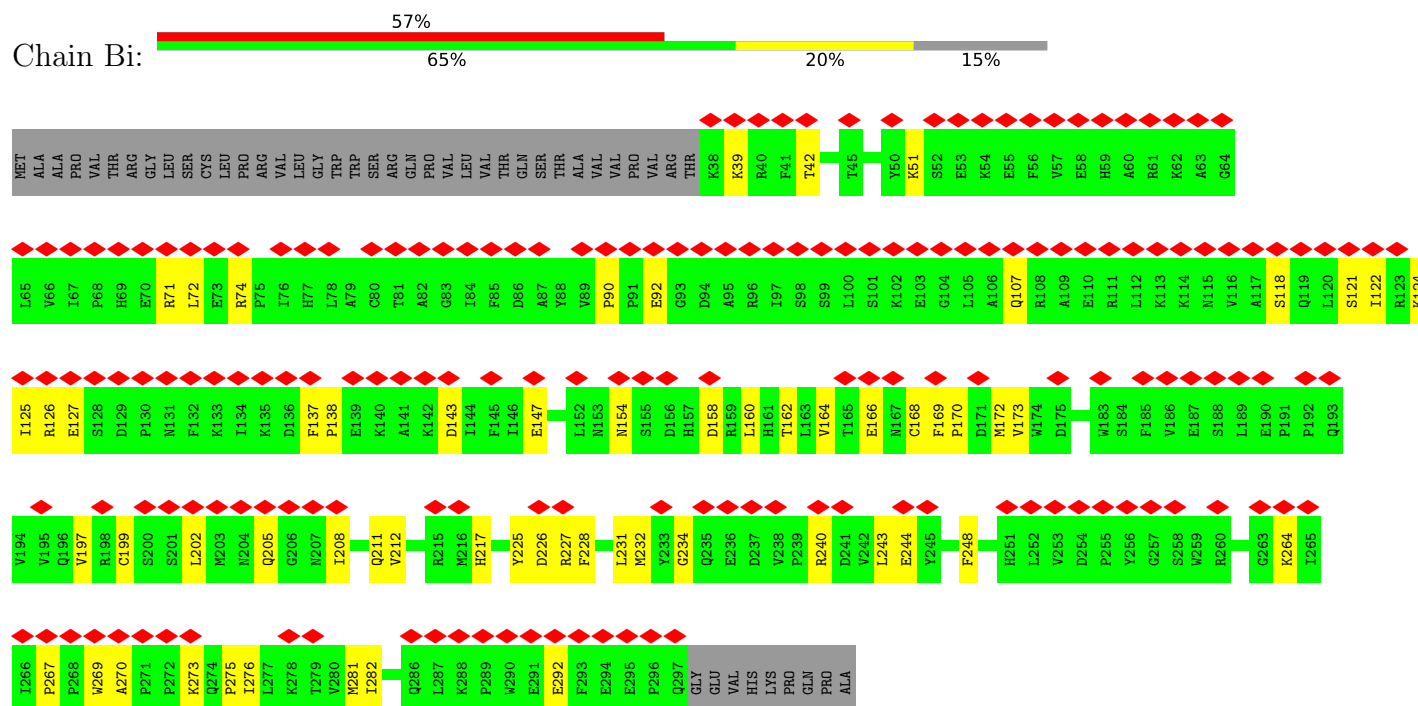
- Molecule 31: Mitochondrial ribosomal protein L43



- Molecule 32: mL44

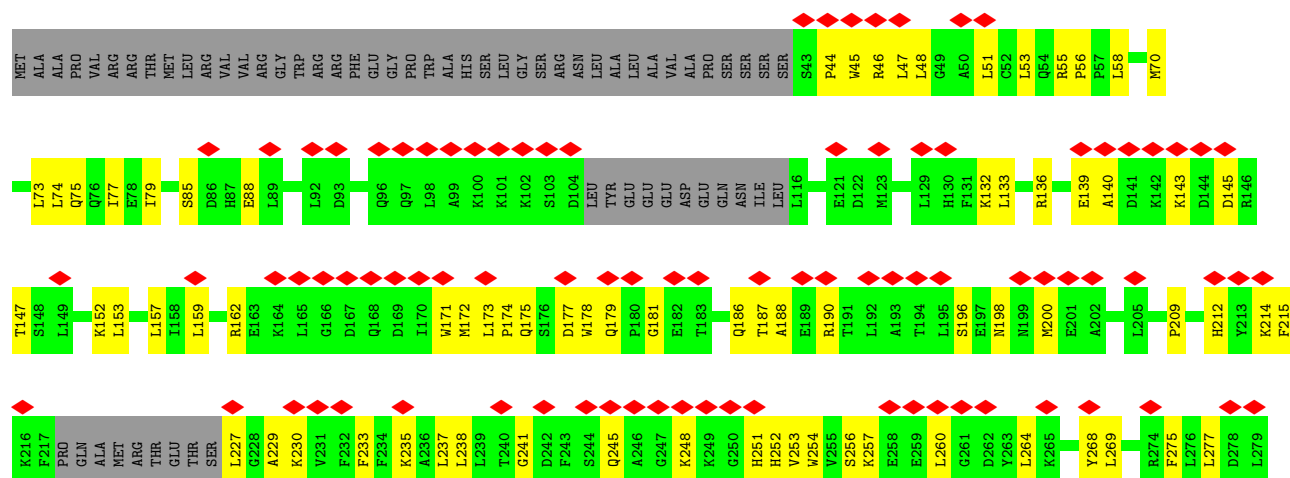


- Molecule 33: Mitochondrial ribosomal protein L45

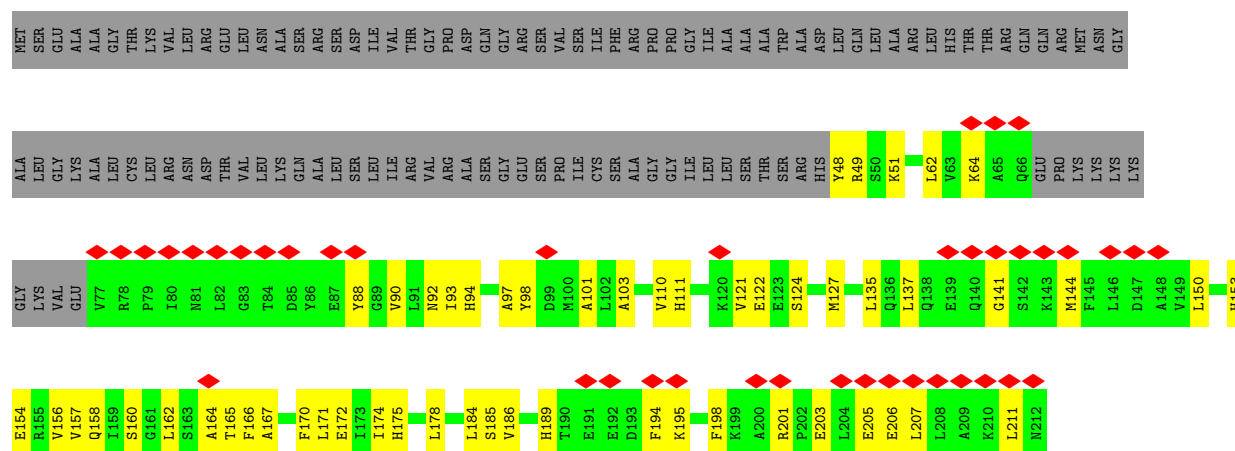


- Molecule 34: Mitochondrial ribosomal protein L46

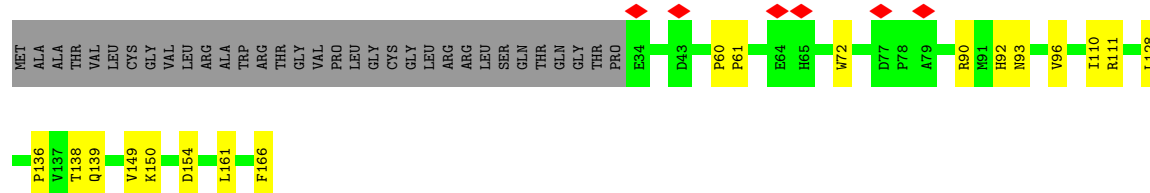




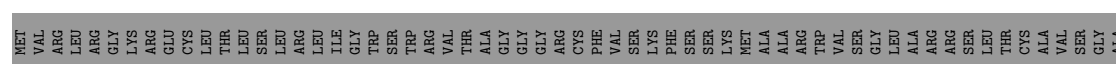
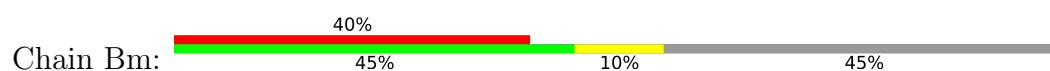
• Molecule 35: 39S ribosomal protein L48, mitochondrial

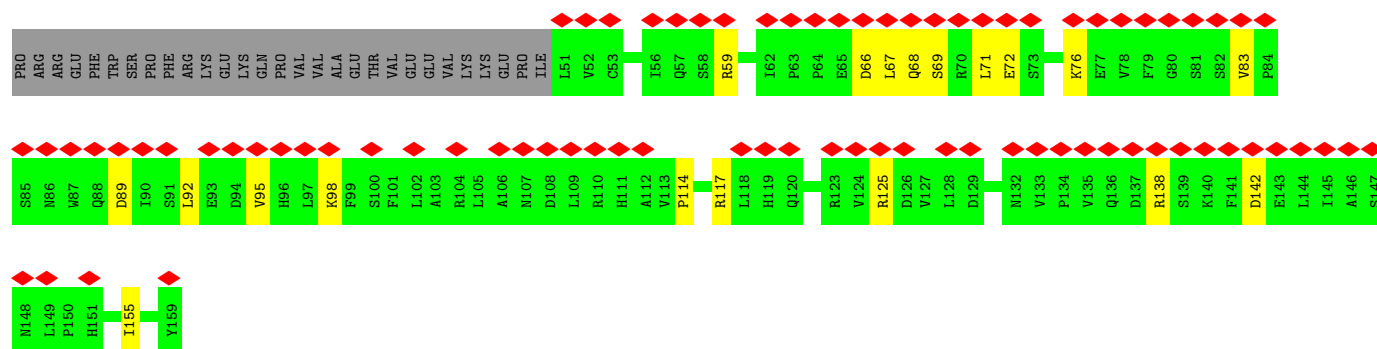


• Molecule 36: Mrpl34

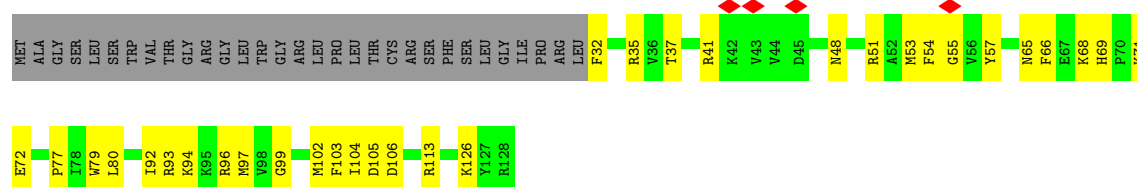


• Molecule 37: Mitochondrial ribosomal protein L50

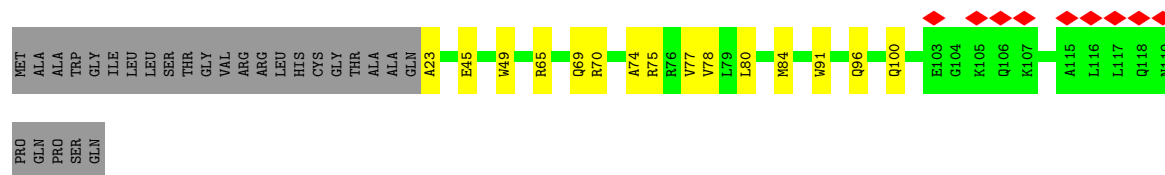




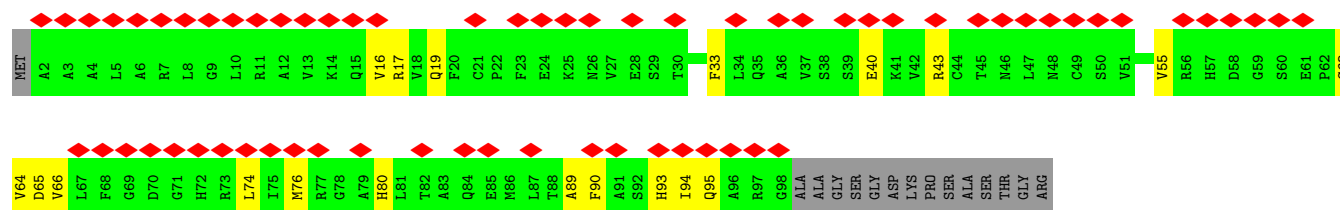
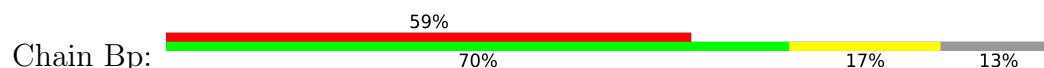
- Molecule 38: Mitochondrial ribosomal protein L51



- Molecule 39: mL52

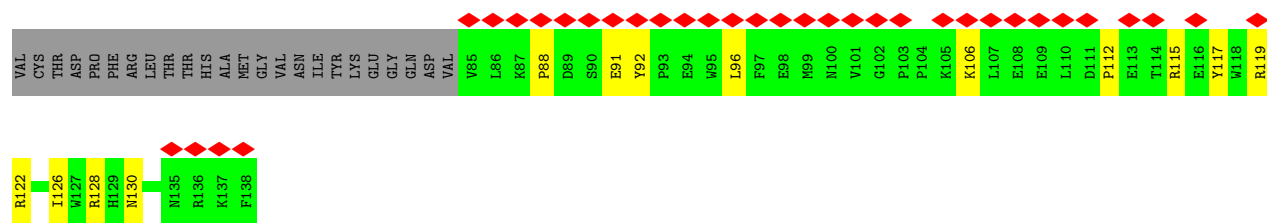


- Molecule 40: mL53



- Molecule 41: mL54

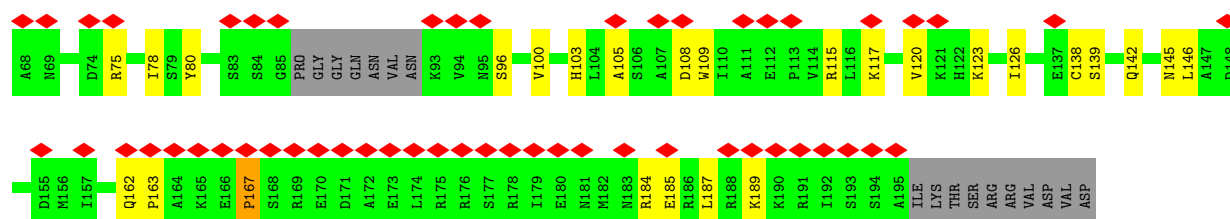
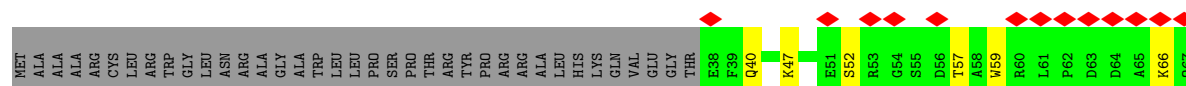




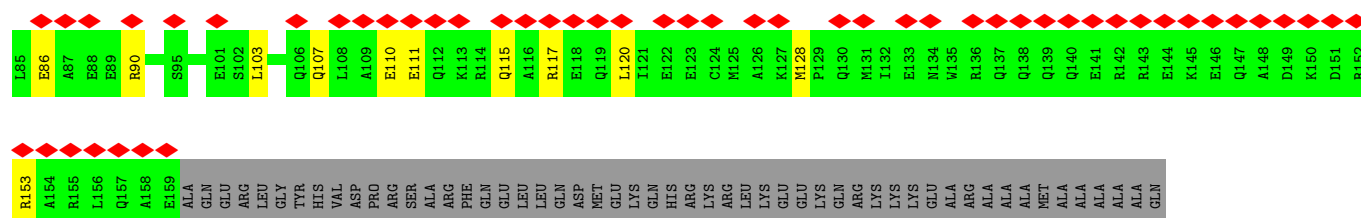
- Molecule 42: Mitochondrial ribosomal protein L57



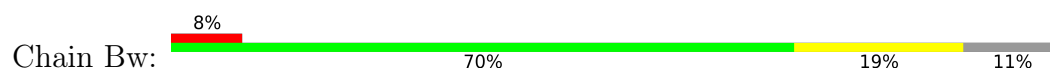
- Molecule 43: mL62 (ICT1)

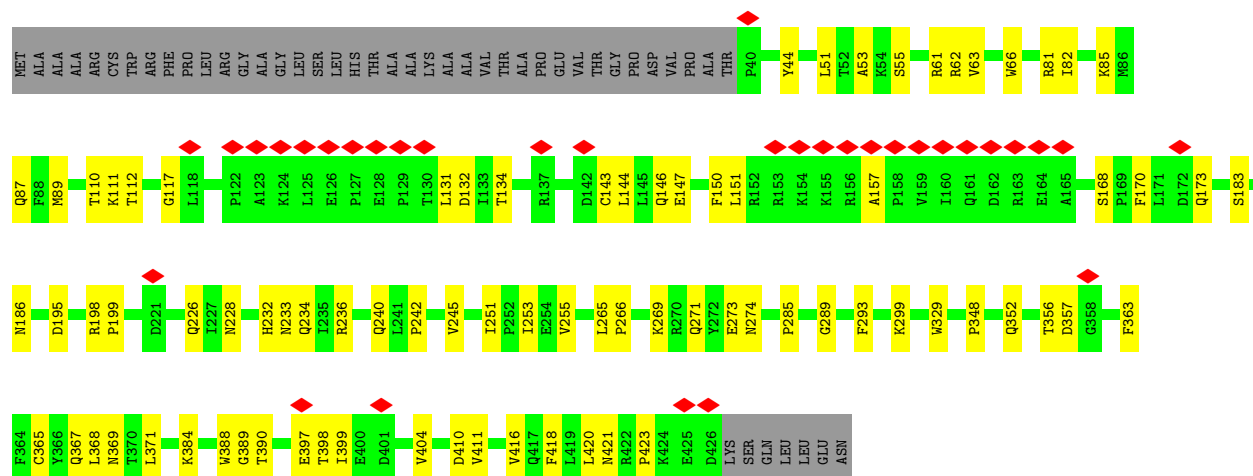


- Molecule 44: mL64

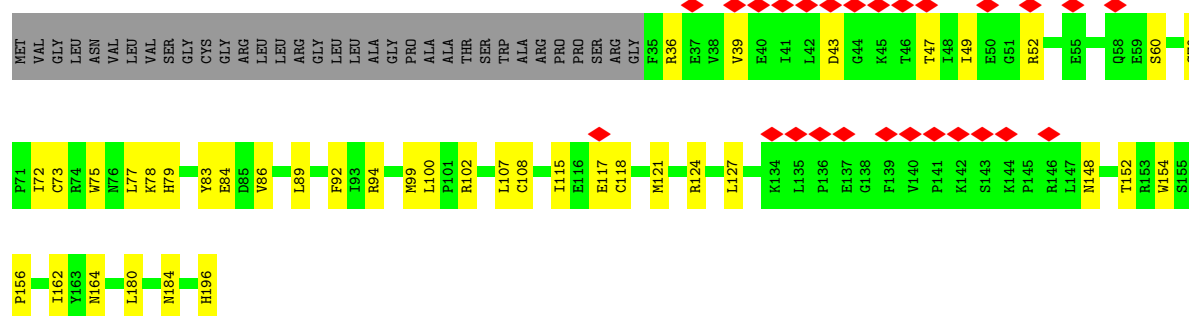


- Molecule 45: 39S ribosomal protein S30, mitochondrial

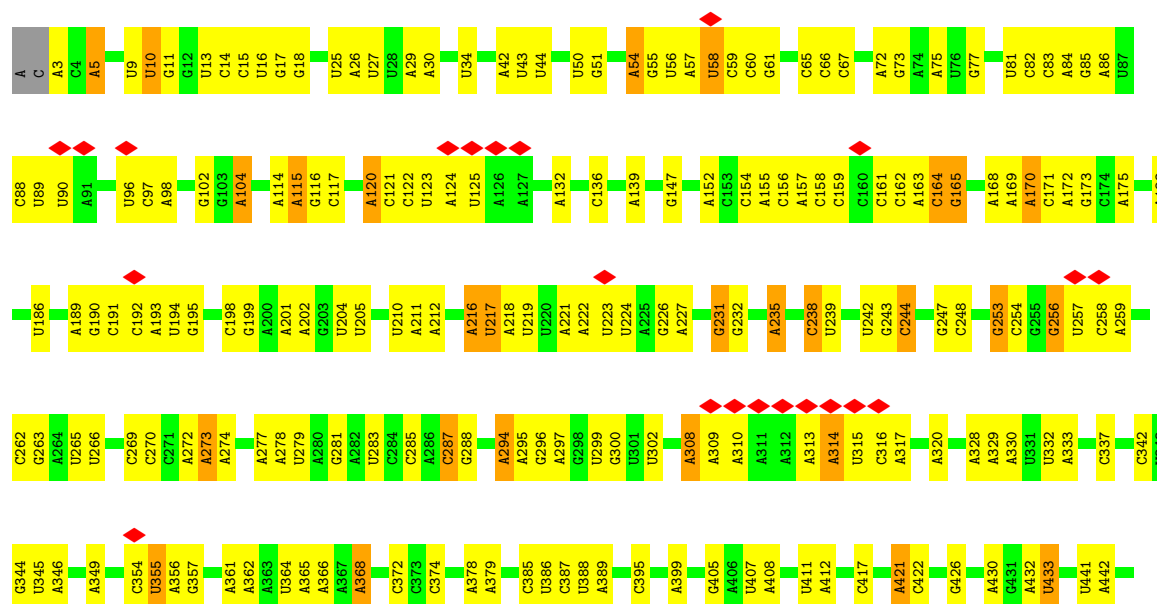


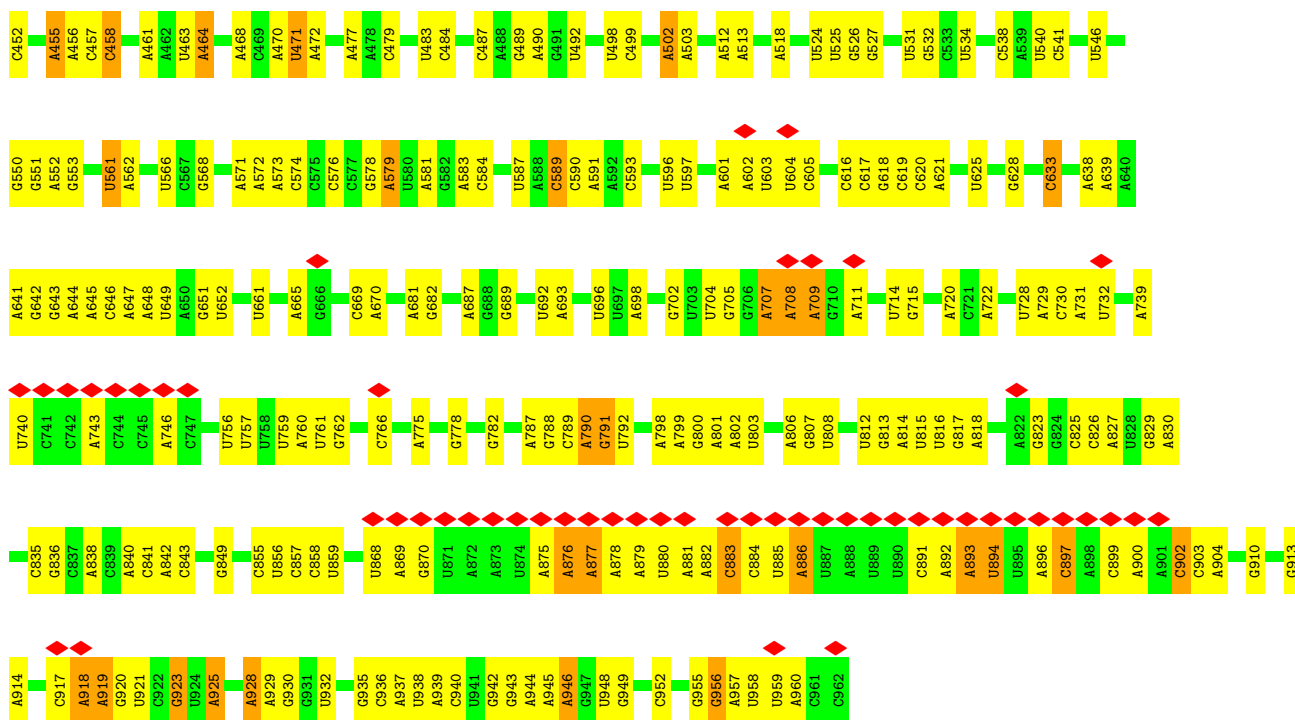


• Molecule 46: Mitochondrial ribosomal protein S18A

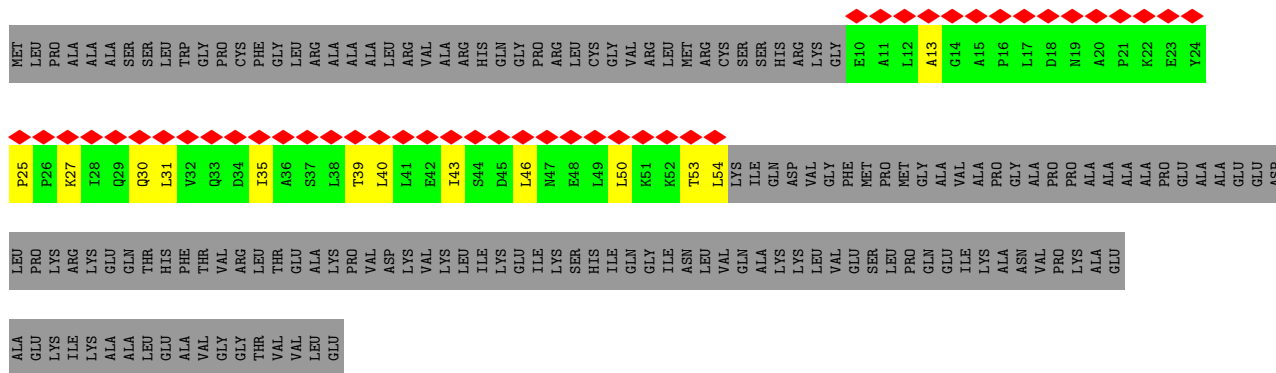


• Molecule 47: 12S rRNA

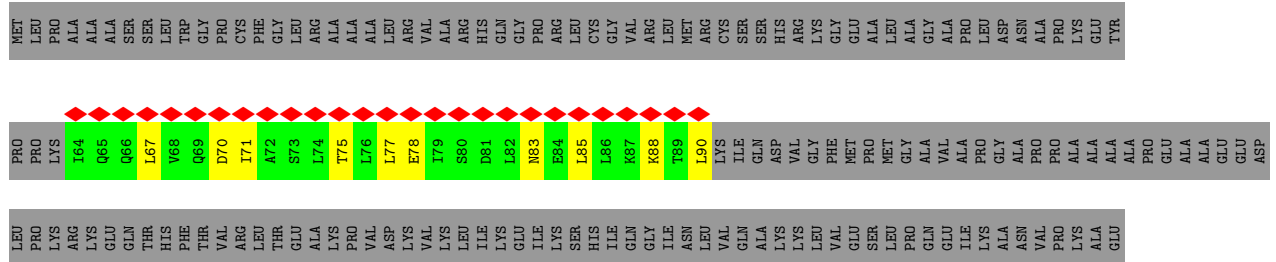


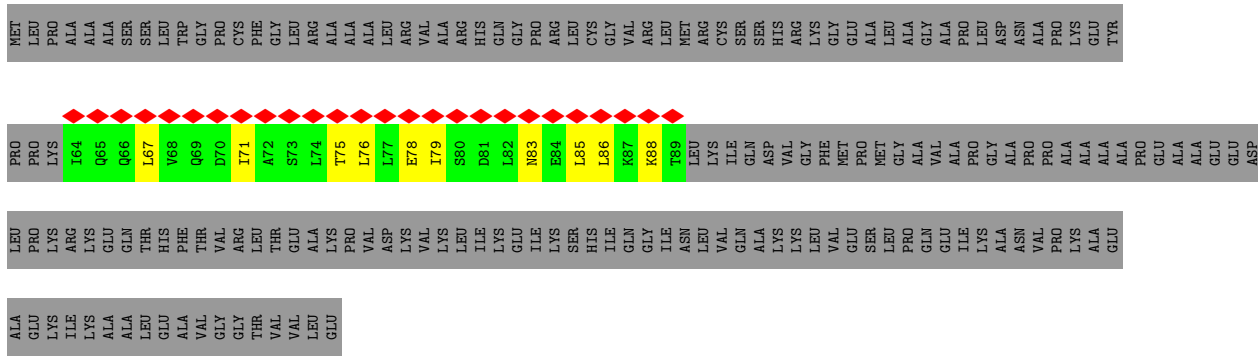


• Molecule 48: Mitochondrial ribosomal protein L12

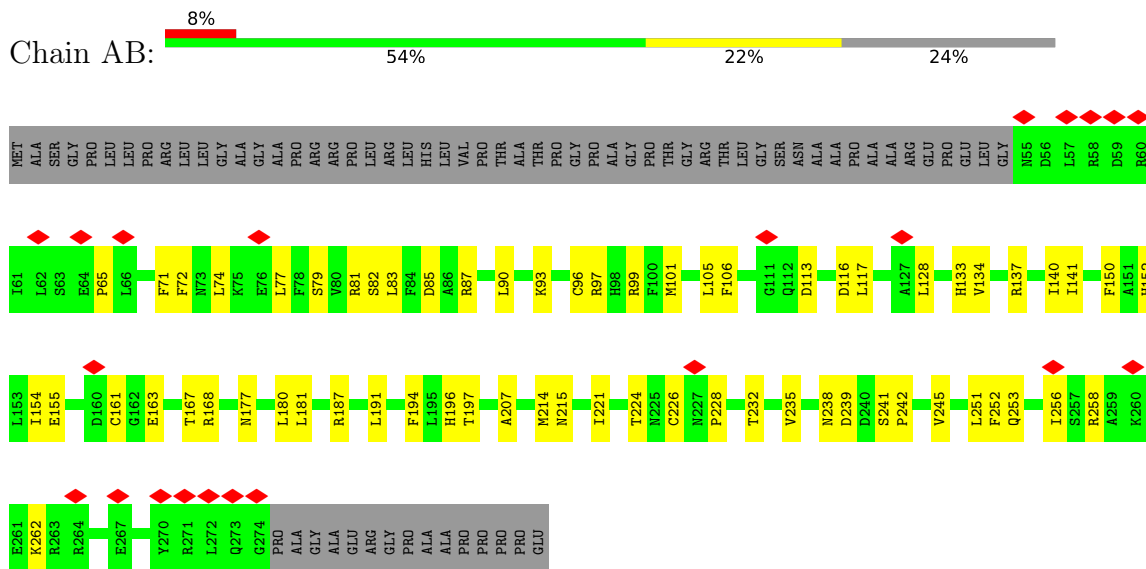


• Molecule 48: Mitochondrial ribosomal protein L12

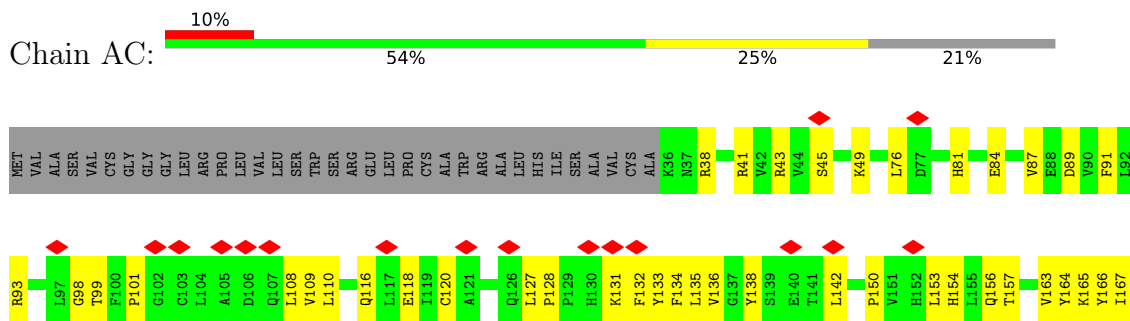




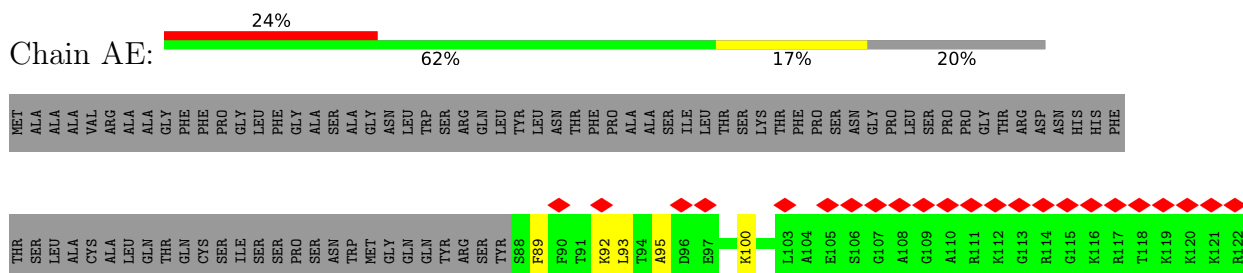
- Molecule 49: Mitochondrial ribosomal protein S2

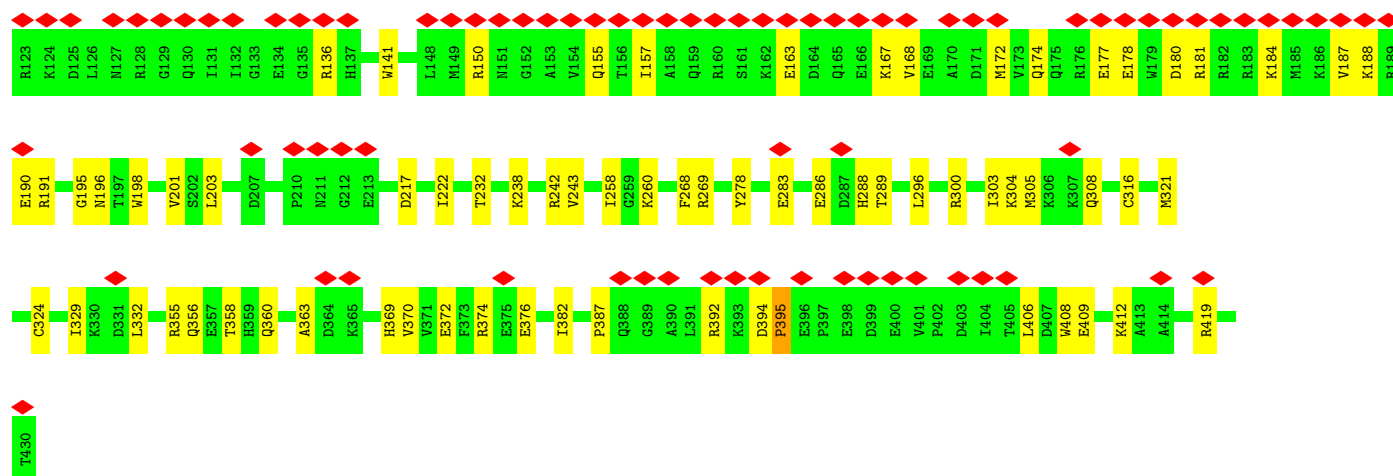


- Molecule 50: Mitochondrial ribosomal protein S24



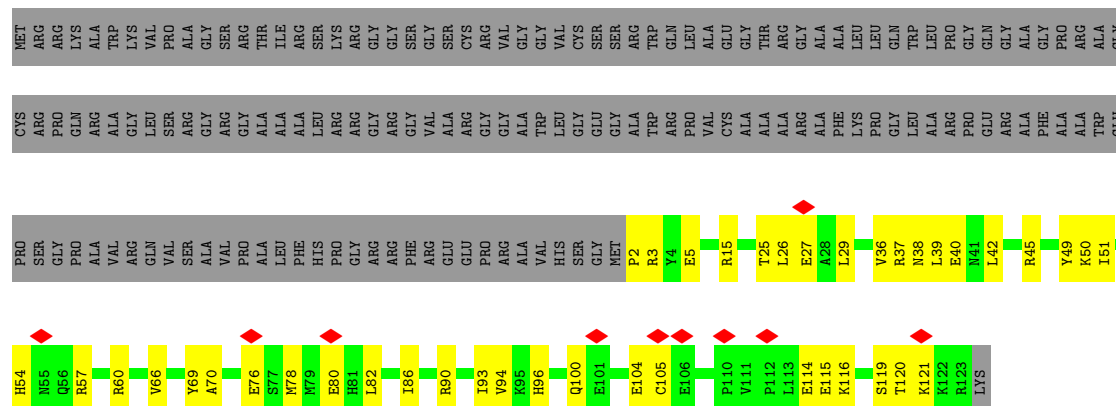
- Molecule 51: Mitochondrial ribosomal protein S5





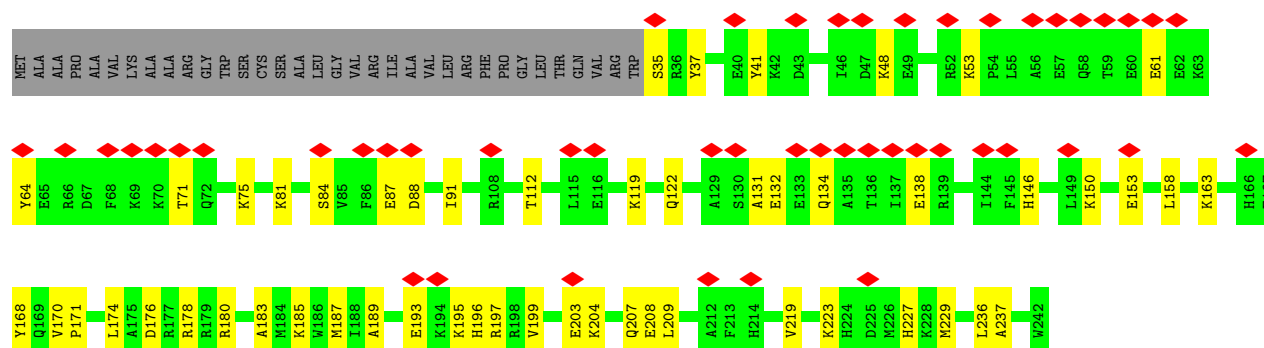
• Molecule 52: bS6m

Chain AF: 29% 15% 56%



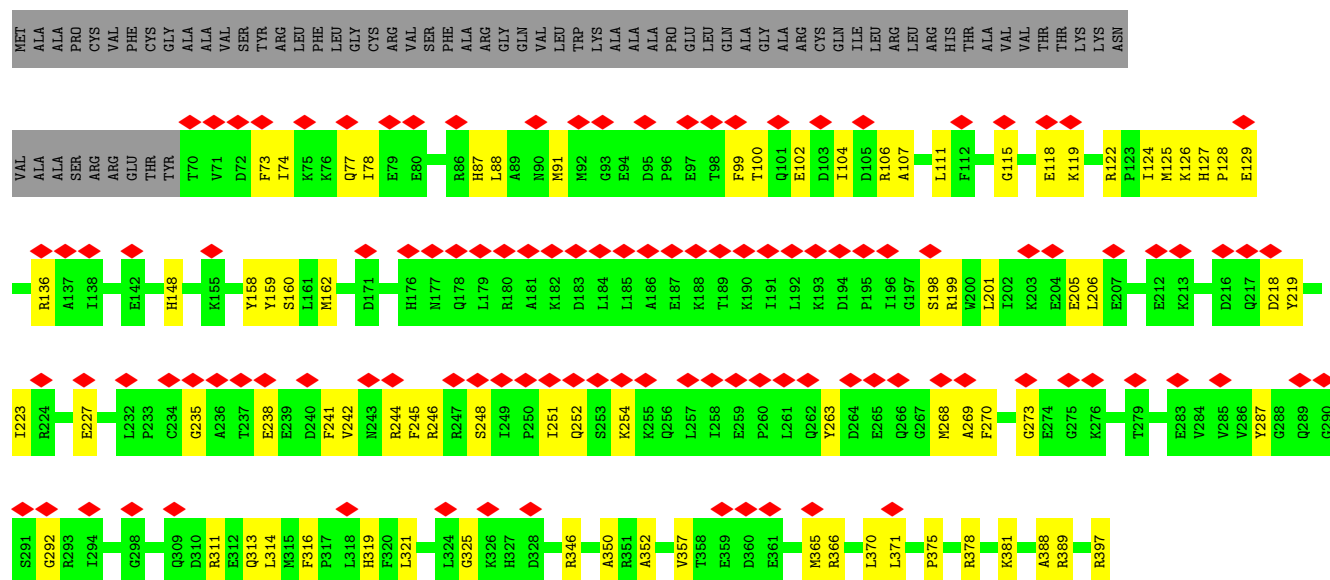
• Molecule 53: Mitochondrial ribosomal protein S7

Chain AG: 20% 64% 22% 14%

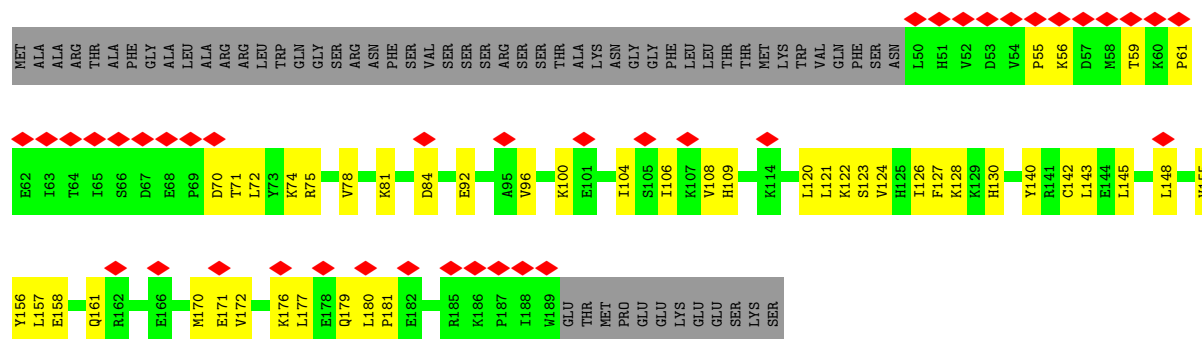


• Molecule 54: uS9m

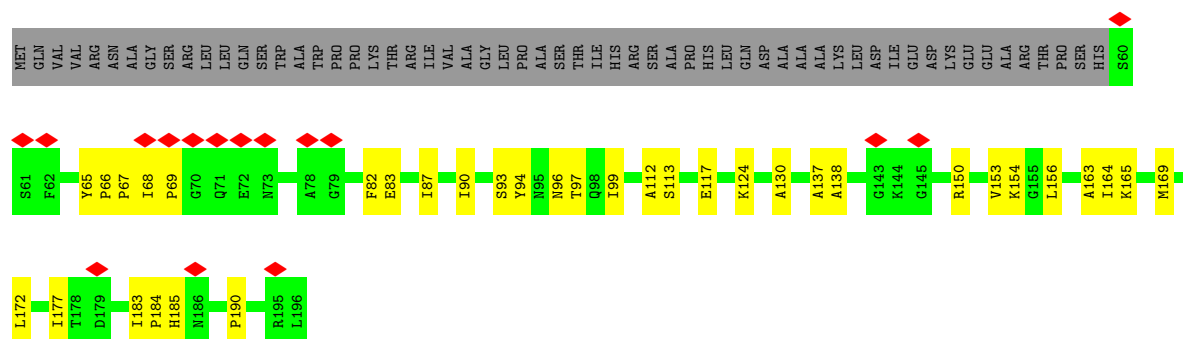
Chain AI: 28% 63% 20% 17%



• Molecule 55: Mitochondrial ribosomal protein S10

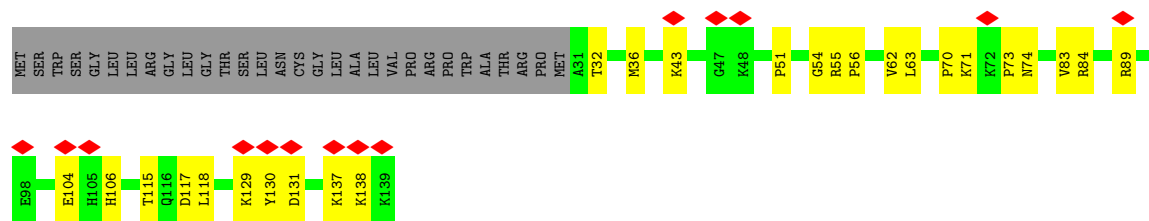


• Molecule 56: uS11m

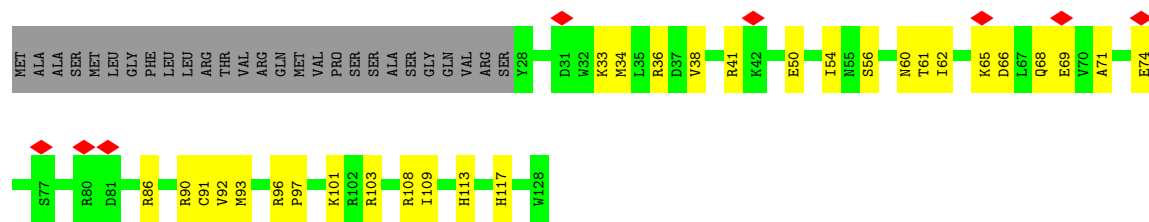


• Molecule 57: Mitochondrial ribosomal protein S12

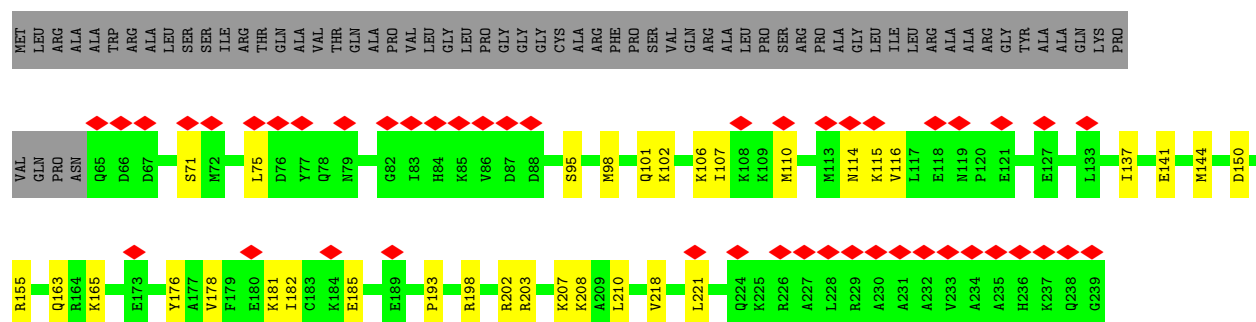




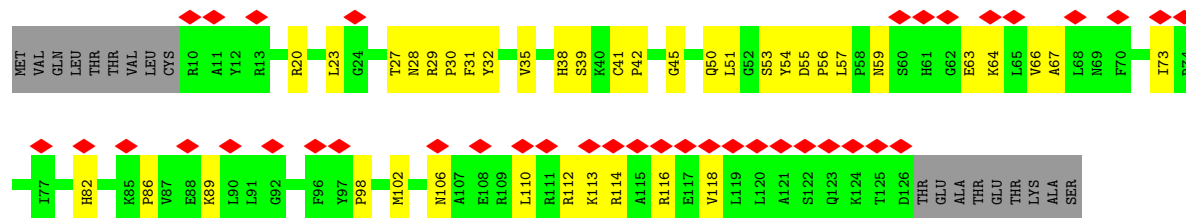
- Molecule 58: Mitochondrial ribosomal protein S14



- Molecule 59: uS15m



- Molecule 60: 28S ribosomal protein S16, mitochondrial

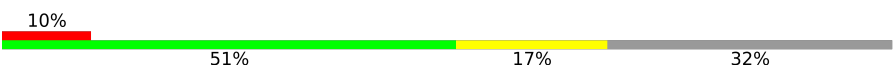


- Molecule 61: uS17m



GLU
ALA
LEU
GLN
LEU
SER
SER
THR
GLN

- Molecule 62: Mitochondrial ribosomal protein S18C

Chain AR: 

MET ALA ALA VAL VAL SER LEU CYS SER ARG LEU GLY ARG ARG LYS SER THR ASP PHE VAL ASN ALA ALA VAL CYS LEU THR ASP PRO GLY THR HIS VAL LEU TRP ARG SER CYS SER GLN TYR LYS GLN VAL ALA S47 M48 E49 D50 L51 P52 V53 P54 M55 E56 M57 P58 Y59 K60

E61 C66 I67 L68 C69 V73 K76 N77 L81 S82 Q83 P87 L100 C101 K104 E107 I117 Y125 K126 D127 P128 L131 V136 Y141 R142 E143

- Molecule 63: bS21m

Chain AU: 

MET A2 N3 H4 A9 R10 T11 Q15 E16 V19 T25 R28 L29 M32 D33 G34 L36 E37 Y45 R52 M66 T72 N73 M76 R80 D82 G86 C87

- Molecule 64: fMet-tRNA^{Met} (P site)

Chain AV: 

A1 G2 U3 A4 A5 G6 G7 U8 C9 A10 A14 A15 A16 U17 A18 C21 U24 C25 C30 C35 A40 A41 A42 A43 U44 G45 U46 U47 G48 G49 U50 U51 A52 U53 A54 C55 C56 C57 U58 U59 C60 C61 C62 G63 U64 A65 C66 U67 A68 C69 C70 A71

- Molecule 65: mRNA

Chain AX: 

A1 A2 A3 G6

- Molecule 66: unknown

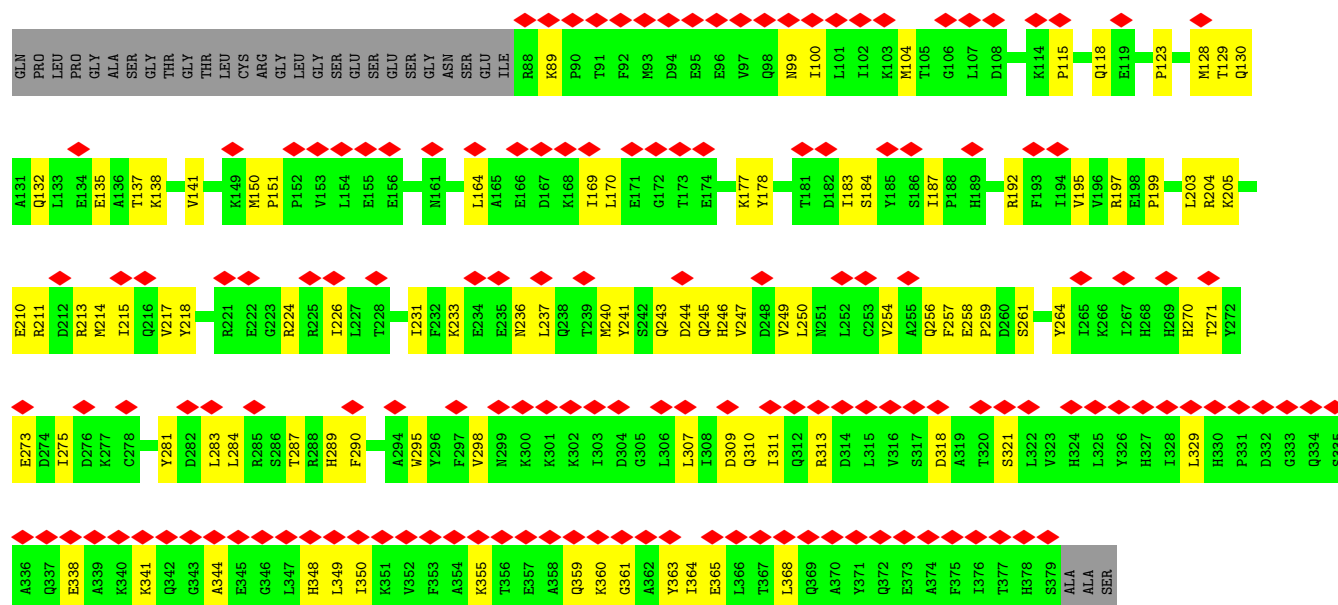
Chain AZ: 

A692 A693 A694 A695 A696 A697 A700 A701 A702 A703 A704 A705 A706 A707 A708 A709

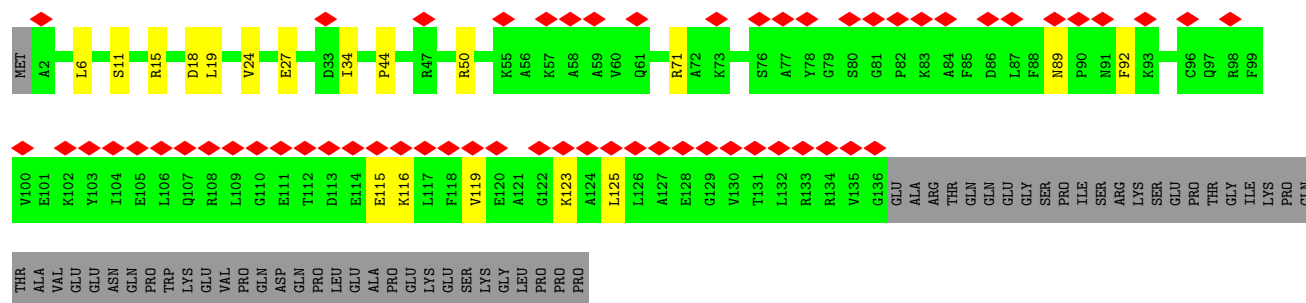
- Molecule 67: Mitochondrial ribosomal protein S22

Chain Aa: 

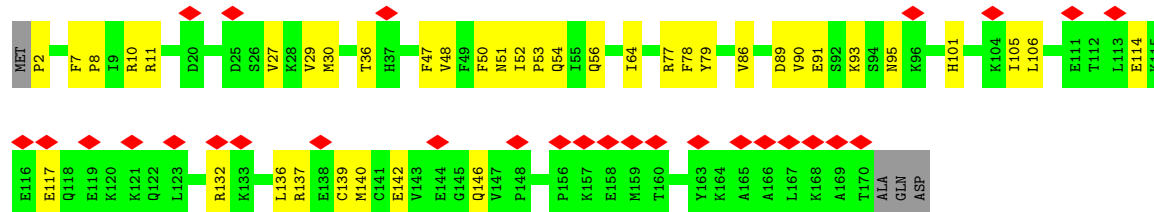
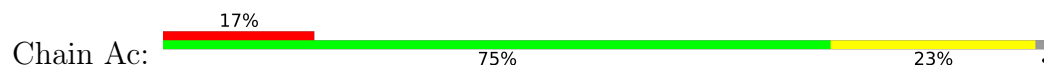
MET PRO HIS GLU PRO ARG ARG GLU THR ARG ASN LEU SER LEU PRO PRO LEU ALA VAL ALA ARG ARG ILE MET ALA THR LEU LYS ALA SER ALA LEU ARG SER LEU GLN THR ASN SER CYS GLY THR GLY ARG VAL CYS PHE PRO VAL ARG ALA ARG PRO ARG PRO ARG ALA LEU LEU



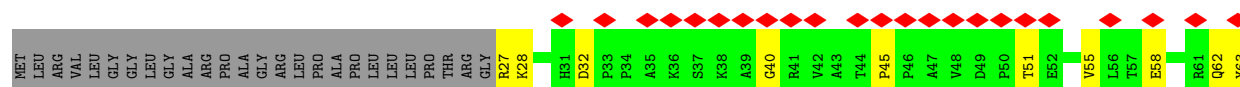
• Molecule 68: mS23

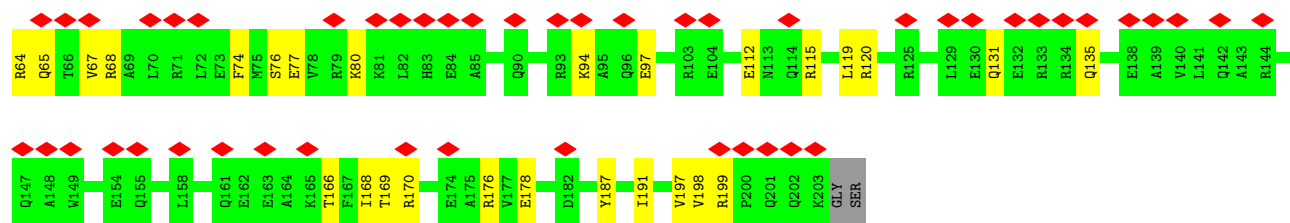


• Molecule 69: Mitochondrial ribosomal protein S25

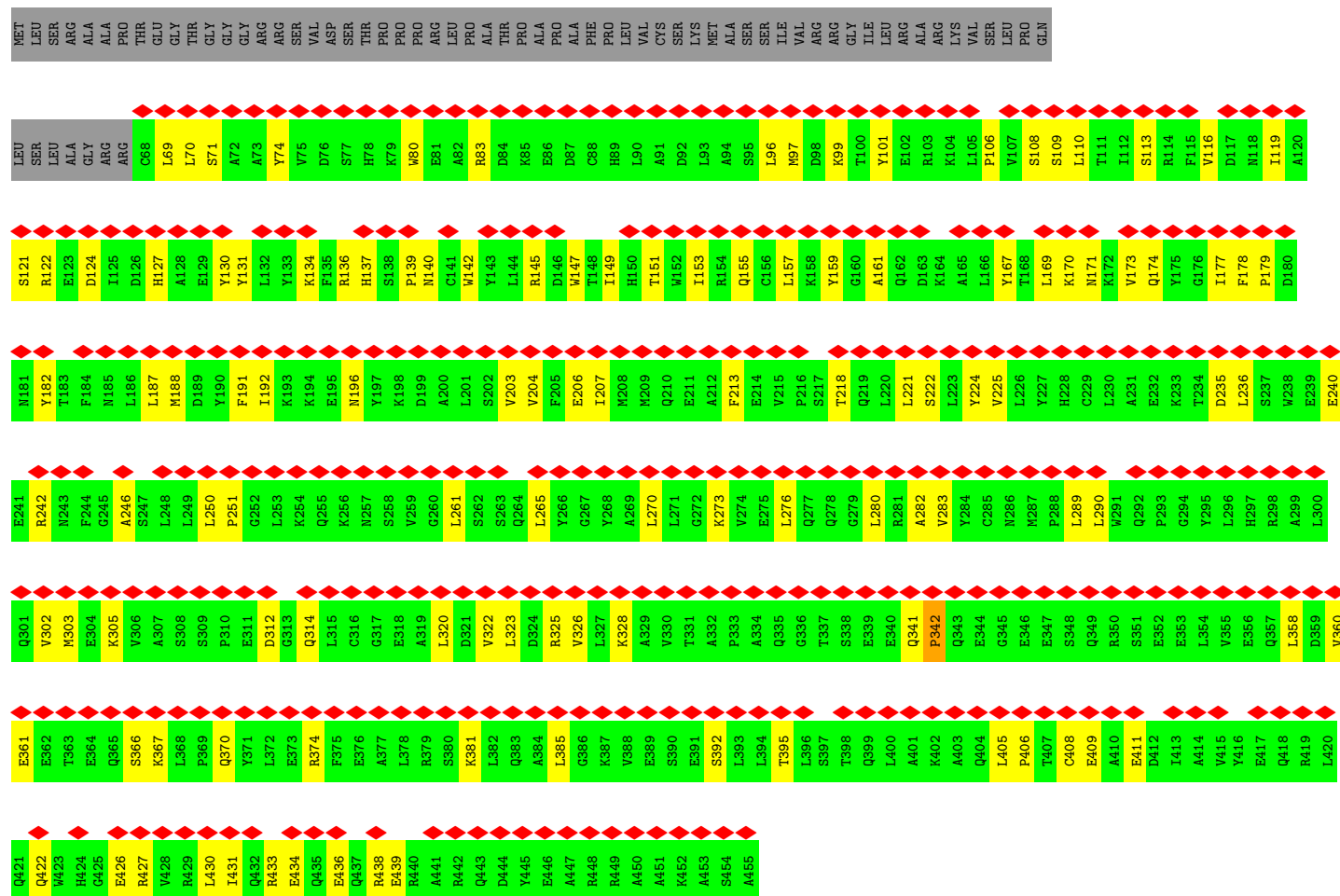
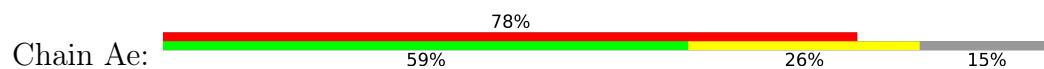


• Molecule 70: Mitochondrial ribosomal protein S26

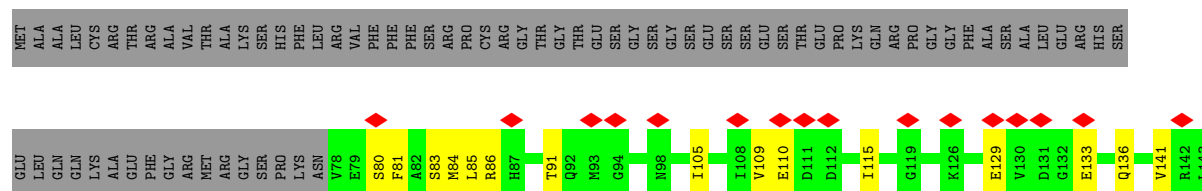


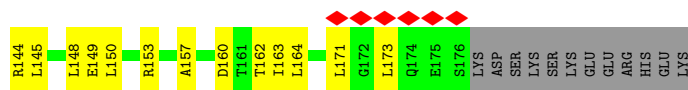


• Molecule 71: Mitochondrial ribosomal protein S27

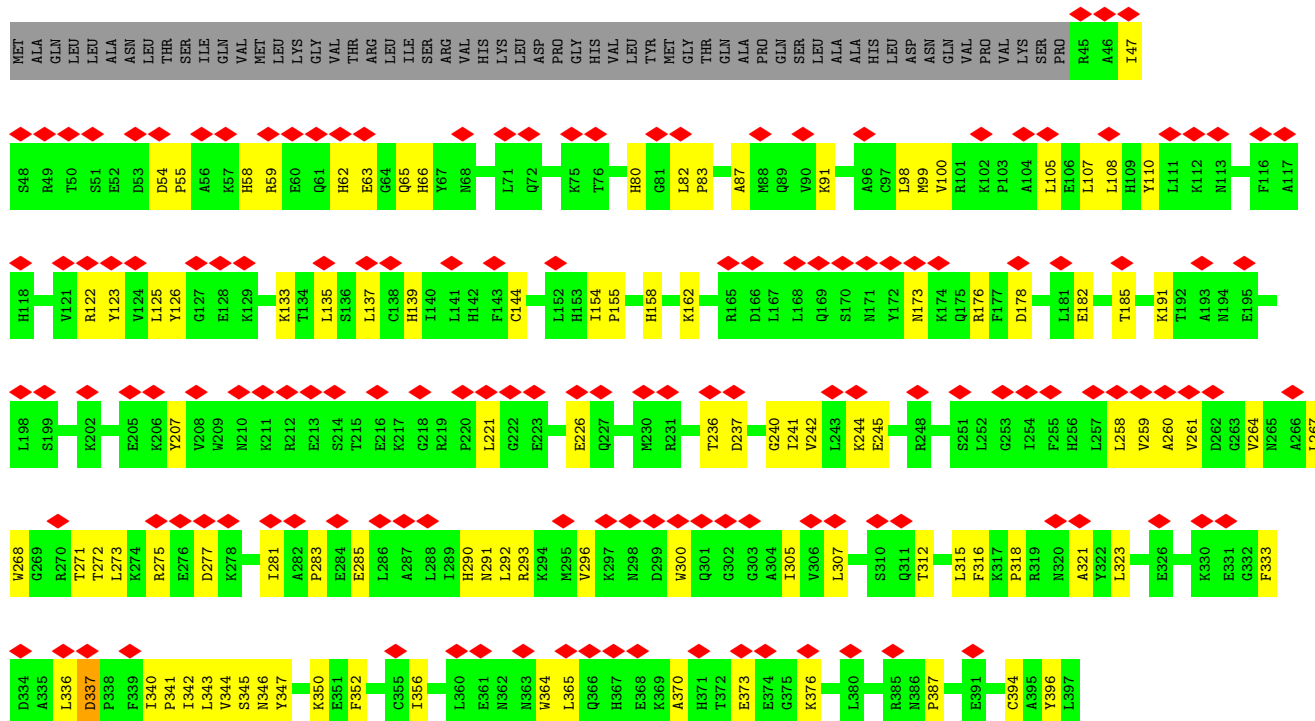
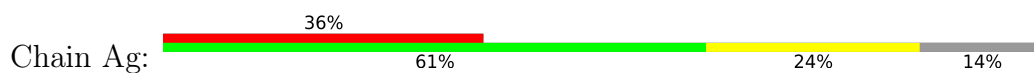


• Molecule 72: Mitoribosomal protein ms28, mrps28

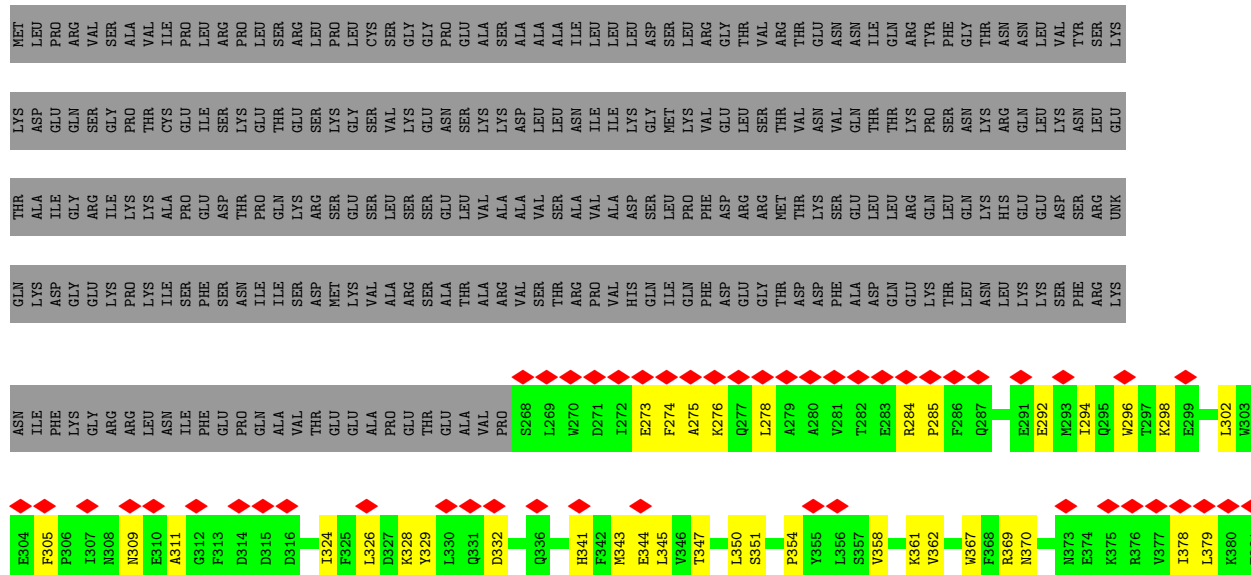


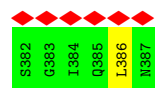


• Molecule 73: Death associated protein 3

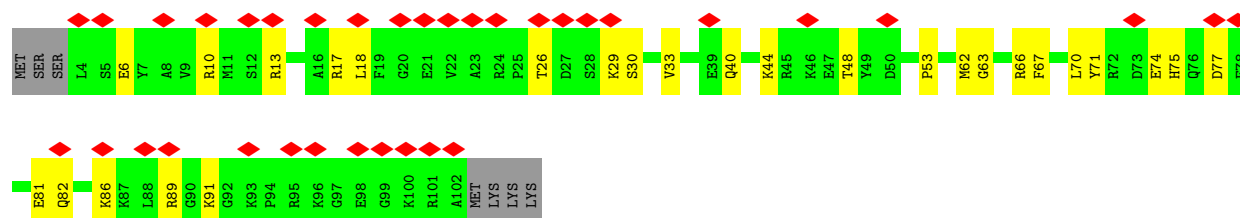


• Molecule 74: Mitochondrial ribosomal protein S31

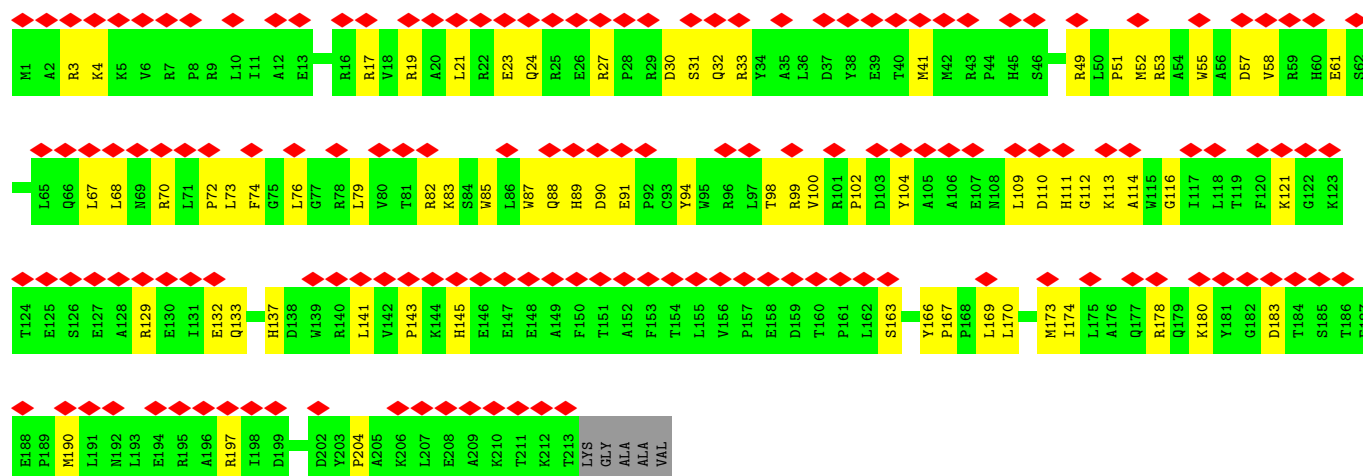




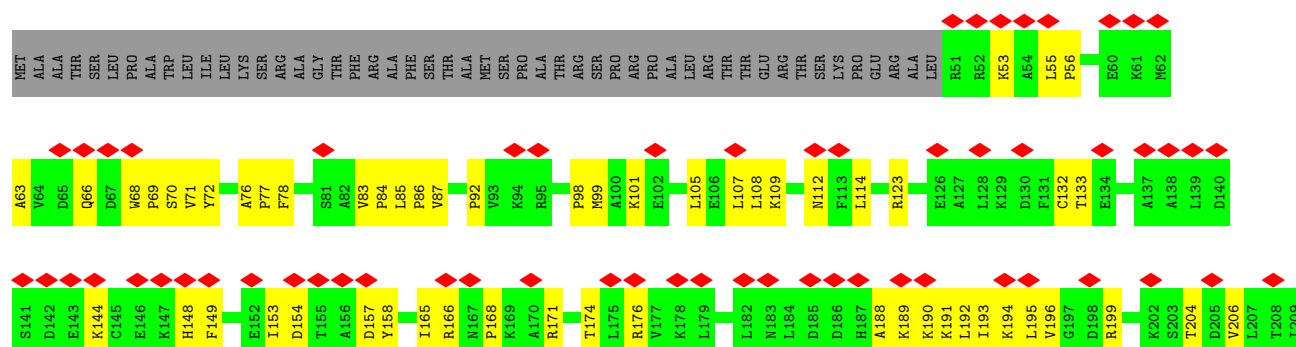
• Molecule 75: mS33

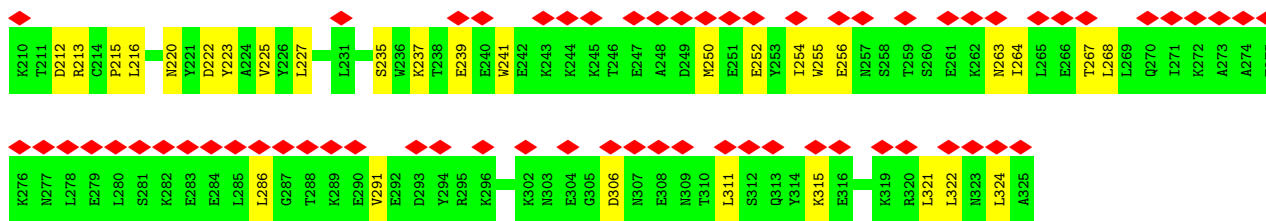


• Molecule 76: mS34

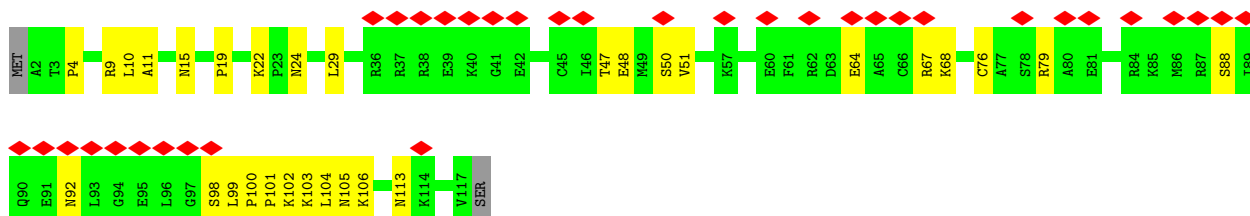
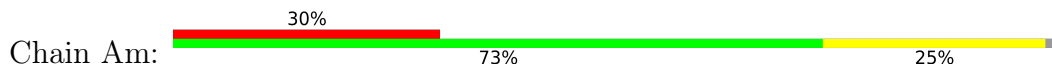


• Molecule 77: Mitochondrial ribosomal protein S35

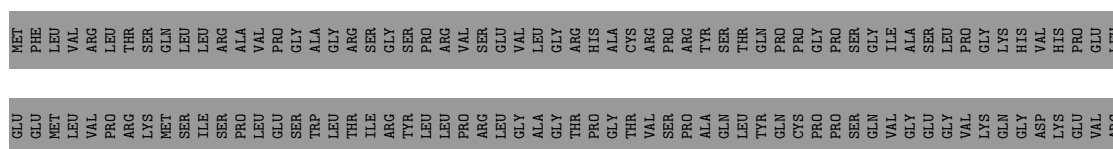




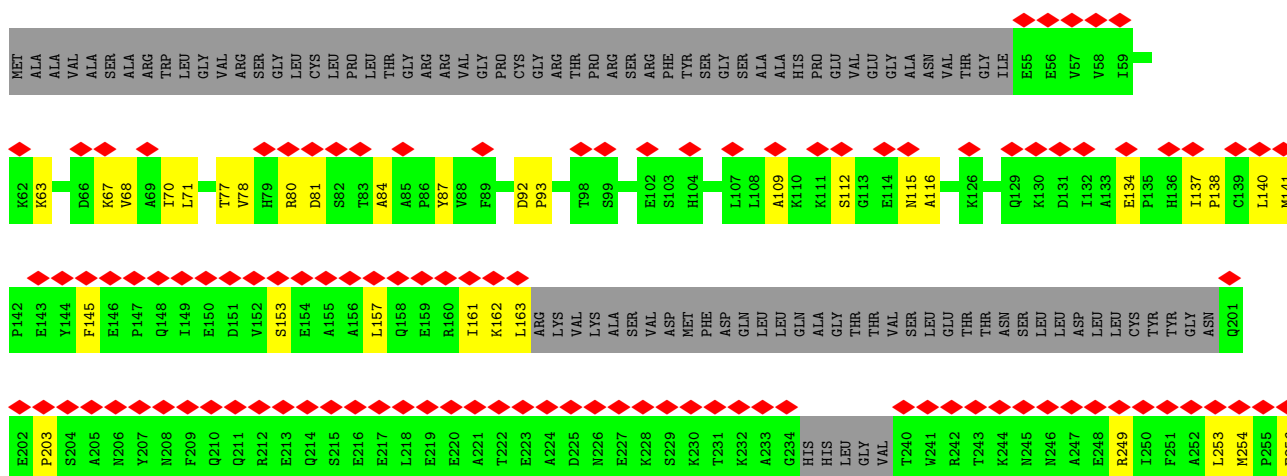
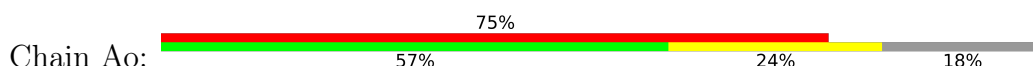
• Molecule 78: mS37

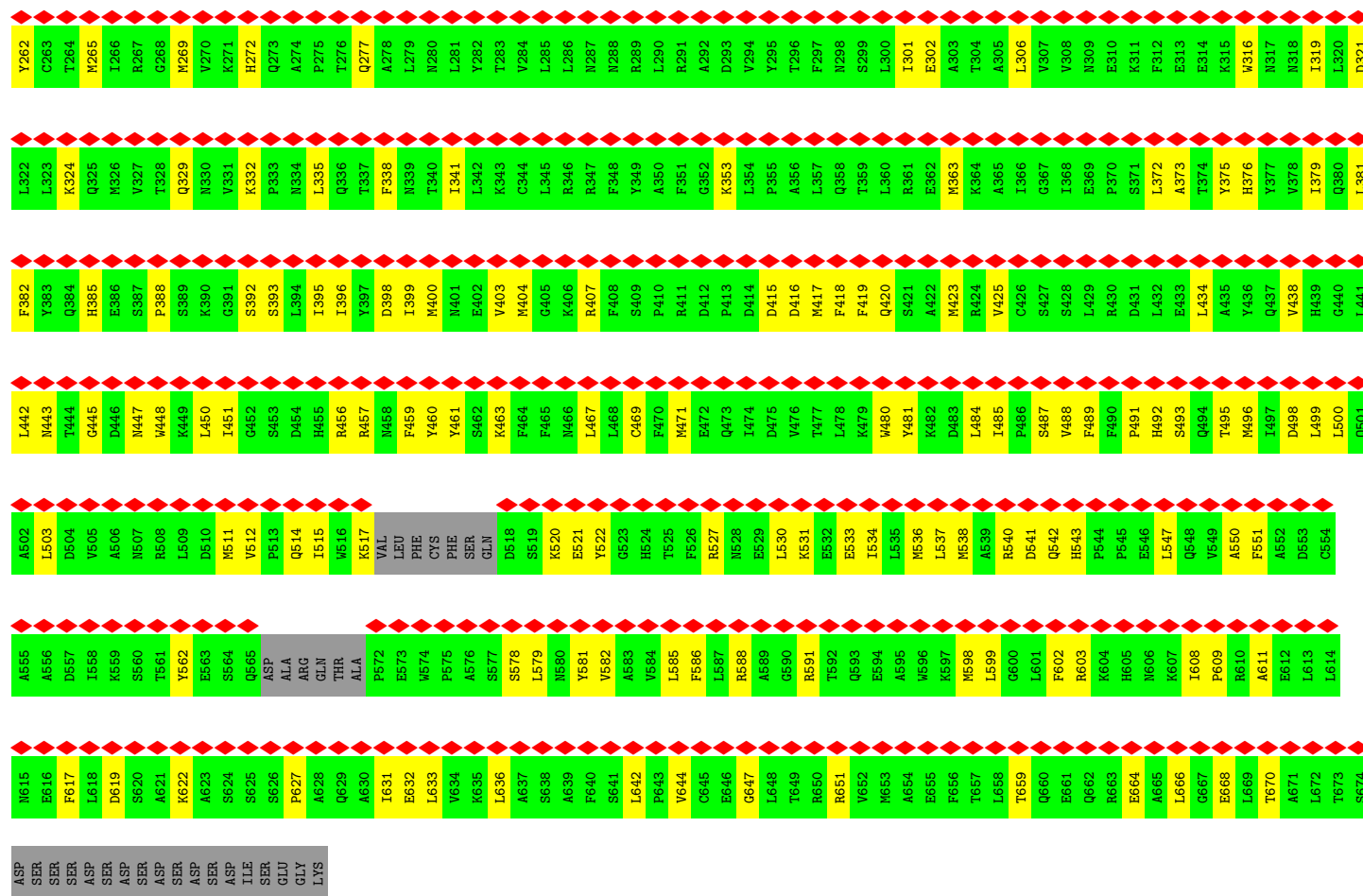


• Molecule 79: Aurora kinase A interacting protein 1



• Molecule 80: Pentatricopeptide repeat domain 3



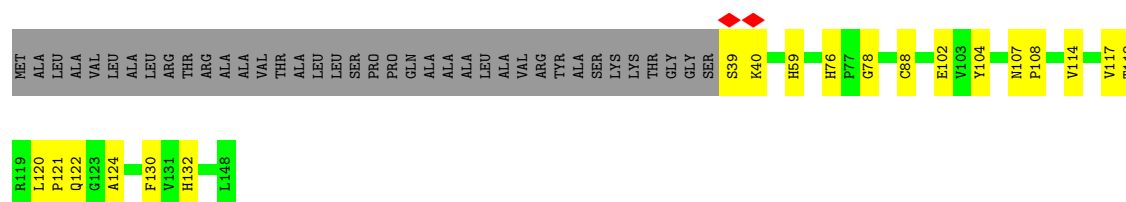


- Molecule 81: 28S ribosomal protein S18b, mitochondrial

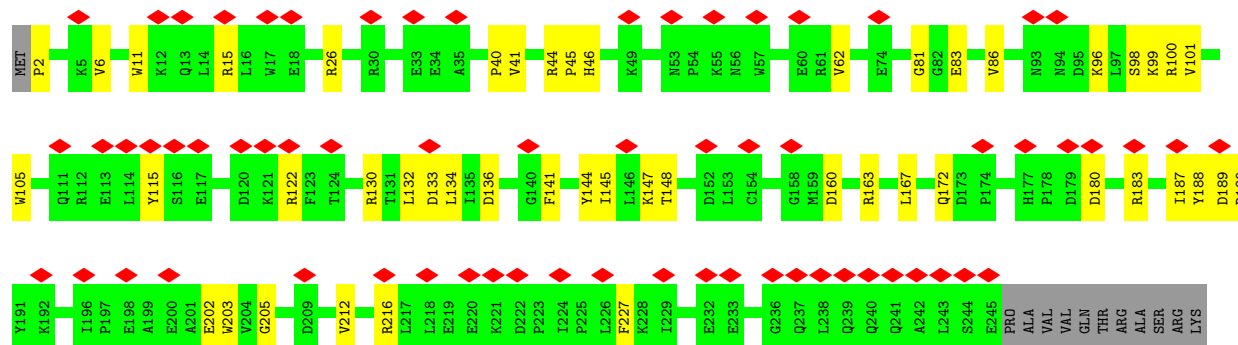
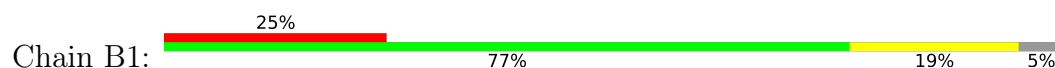


- Molecule 82: Mitochondrial ribosomal protein L27

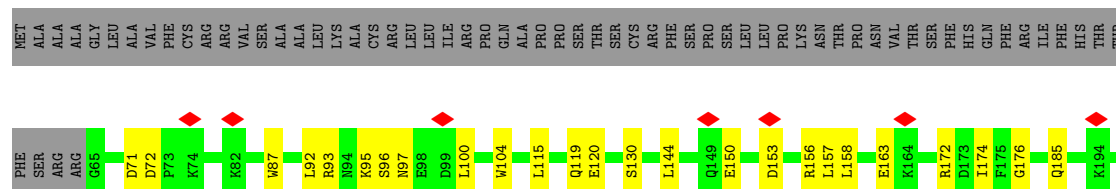




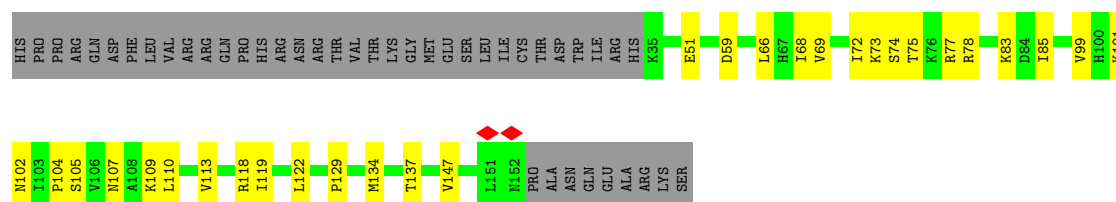
• Molecule 83: Mitochondrial ribosomal protein L28



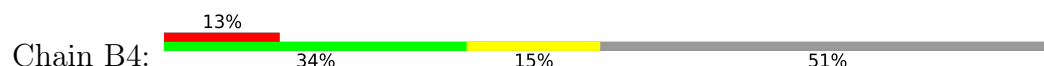
• Molecule 84: Mitochondrial ribosomal protein L47

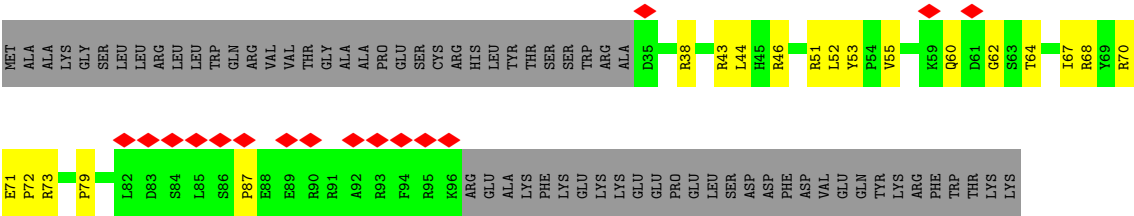


• Molecule 85: uL30m

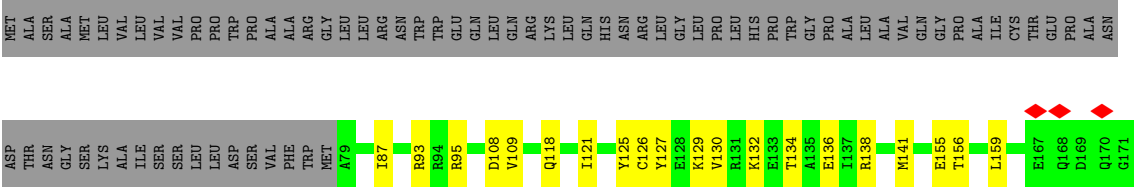


• Molecule 86: bL31m

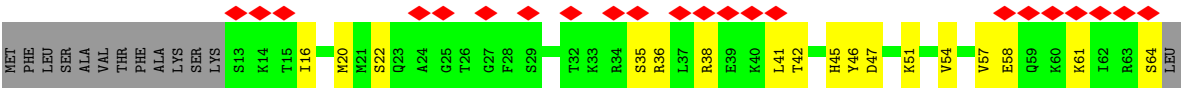




• Molecule 87: bL32m



• Molecule 88: bL33m



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	93623	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	40	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	FEI FALCON III (4k x 4k)	Depositor
Maximum map value	1.037	Depositor
Minimum map value	-0.651	Depositor
Average map value	0.001	Depositor
Map value standard deviation	0.046	Depositor
Recommended contour level	0.13	Depositor
Map size (Å)	444.8, 444.8, 444.8	wwPDB
Map dimensions	320, 320, 320	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.39, 1.39, 1.39	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: GTP, MG, SPM, ZN, 5GP

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	B7	0.09	0/395	0.23	0/524
2	B8	0.09	0/853	0.24	0/1136
3	B9	0.10	0/342	0.27	0/450
4	BA	0.10	0/36784	0.24	0/57270
5	BB	0.08	0/1595	0.20	0/2475
6	BD	0.12	0/1898	0.35	0/2555
7	BE	0.12	0/2493	0.33	0/3387
8	BF	0.12	0/2069	0.35	0/2816
9	BI	0.10	0/819	0.31	0/1101
10	BJ	0.13	0/1742	0.36	0/2358
11	BK	0.11	0/1323	0.32	0/1785
12	BL	0.10	0/1127	0.28	0/1512
13	BN	0.11	0/1487	0.28	0/2017
14	BO	0.20	0/912	0.38	0/1231
15	BP	0.11	0/2368	0.30	0/3198
16	BQ	0.11	0/1850	0.28	0/2491
17	BR	0.11	0/1262	0.30	0/1700
18	BS	0.12	0/1197	0.36	0/1624
19	BT	0.17	0/2002	0.35	1/2708 (0.0%)
20	BU	0.13	0/1179	0.31	0/1578
21	BV	0.12	0/1256	0.29	0/1706
22	BW	0.12	0/1407	0.32	0/1891
23	BX	0.13	0/1211	0.29	0/1646
24	BY	0.10	0/1719	0.32	0/2329
25	Ba	0.11	0/3267	0.30	0/4455
26	Bb	0.12	0/3047	0.29	0/4139
27	Bc	0.13	0/2464	0.35	0/3330
28	Bd	0.15	0/1183	0.36	0/1594
29	Be	0.11	0/1000	0.27	0/1345
30	Bf	0.14	0/851	0.50	2/1159 (0.2%)
31	Bg	0.14	0/1191	0.37	0/1614
32	Bh	0.11	0/2372	0.31	0/3211

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
33	Bi	0.12	0/2199	0.33	0/2980
34	Bj	0.12	0/1811	0.37	0/2436
35	Bk	0.15	0/1270	0.40	0/1714
36	Bl	0.13	0/1135	0.31	0/1549
37	Bm	0.12	0/917	0.31	0/1248
38	Bn	0.12	0/860	0.33	0/1150
39	Bo	0.10	0/787	0.26	0/1056
40	Bp	0.12	0/752	0.38	0/1013
41	Bq	0.13	0/490	0.33	0/662
42	Bt	0.11	0/798	0.28	0/1073
43	Bu	0.10	0/1214	0.34	1/1630 (0.1%)
44	Bv	0.09	0/1157	0.23	0/1560
45	Bw	0.12	0/3206	0.31	0/4354
46	Bx	0.14	0/1364	0.34	0/1849
47	AA	0.09	0/22852	0.21	0/35580
48	CL	0.11	0/319	0.26	0/435
48	DL	0.10	0/212	0.28	0/286
48	EL	0.10	0/221	0.28	0/297
48	FL	0.10	0/212	0.27	0/286
48	GL	0.09	0/212	0.28	0/286
48	HL	0.10	0/204	0.28	0/275
49	AB	0.15	0/1804	0.40	0/2445
50	AC	0.13	0/1105	0.35	0/1496
51	AE	0.20	0/2785	0.41	2/3735 (0.1%)
52	AF	0.13	0/999	0.36	0/1347
53	AG	0.13	0/1763	0.36	0/2368
54	AI	0.17	0/2707	0.34	0/3636
55	AJ	0.15	0/1181	0.36	0/1597
56	AK	0.12	0/1027	0.33	0/1389
57	AL	0.12	0/858	0.32	0/1152
58	AN	0.11	0/874	0.33	0/1171
59	AO	0.11	0/1473	0.34	0/1970
60	AP	0.15	0/954	0.37	0/1284
61	AQ	0.10	0/894	0.27	0/1213
62	AR	0.12	0/802	0.37	0/1079
63	AU	0.10	0/745	0.31	0/993
64	AV	0.10	0/1673	0.26	0/2602
65	AX	0.13	0/147	0.38	0/227
66	AZ	0.08	0/89	0.31	0/123
67	Aa	0.13	0/2428	0.42	2/3279 (0.1%)
68	Ab	0.12	0/1126	0.32	0/1514
69	Ac	0.13	0/1399	0.38	0/1881
70	Ad	0.14	0/1490	0.36	0/2005

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
71	Ae	0.14	0/3171	0.44	1/4292 (0.0%)
72	Af	0.12	0/790	0.35	0/1064
73	Ag	0.13	0/2945	0.37	0/3984
74	Ah	0.12	0/1045	0.33	0/1409
75	Ai	0.13	0/841	0.44	0/1121
76	Aj	0.12	0/1835	0.34	0/2484
77	Ak	0.13	0/2268	0.37	0/3069
78	Am	0.12	0/947	0.36	0/1268
79	An	0.10	0/650	0.32	0/858
80	Ao	0.14	0/4626	0.40	1/6269 (0.0%)
81	Ap	0.13	0/1616	0.38	0/2195
82	B0	0.11	0/880	0.28	0/1189
83	B1	0.10	0/2093	0.27	0/2835
84	B2	0.10	0/1586	0.30	0/2123
85	B3	0.13	0/993	0.33	0/1341
86	B4	0.14	0/481	0.48	1/653 (0.2%)
87	B5	0.10	0/917	0.27	0/1227
88	B6	0.14	0/430	0.41	0/570
All	All	0.12	0/183294	0.31	11/260511 (0.0%)

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
6	BD	0	1
10	BJ	0	1
38	Bn	0	1
73	Ag	0	1
All	All	0	4

There are no bond length outliers.

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
51	AE	395	PRO	CA-N-CD	-9.30	98.97	112.00
67	Aa	257	PHE	CA-C-N	7.62	135.43	120.94
67	Aa	257	PHE	C-N-CA	7.62	135.43	120.94
30	Bf	80	PRO	N-CA-CB	7.54	111.16	103.25
71	Ae	342	PRO	N-CA-CB	7.48	111.11	103.25
30	Bf	78	PRO	N-CA-CB	7.36	110.98	103.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
43	Bu	167	PRO	N-CA-CB	7.34	110.96	103.25
51	AE	395	PRO	N-CD-CG	-6.91	92.83	103.20
86	B4	87	PRO	N-CA-CB	6.34	110.02	103.23
19	BT	68	PRO	N-CA-CB	6.34	109.91	103.25
80	Ao	203	PRO	N-CA-CB	6.23	110.11	103.39

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
73	Ag	337	ASP	Peptide
6	BD	207	TYR	Peptide
10	BJ	64	CYS	Peptide
38	Bn	65	ASN	Peptide

5.2 Too-close contacts [i](#)

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	B7	387	0	413	9	0
2	B8	833	0	883	14	0
3	B9	335	0	359	4	0
4	BA	32844	0	16616	481	0
5	BB	1427	0	726	17	0
6	BD	1860	0	1923	37	0
7	BE	2420	0	2418	60	0
8	BF	2011	0	2049	71	0
9	BI	805	0	845	15	0
10	BJ	1705	0	1804	58	0
11	BK	1303	0	1343	35	0
12	BL	1113	0	1148	32	0
13	BN	1444	0	1437	35	0
14	BO	896	0	946	21	0
15	BP	2312	0	2373	59	0
16	BQ	1803	0	1841	37	0
17	BR	1240	0	1260	32	0
18	BS	1168	0	1159	32	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
19	BT	1954	0	1943	34	0
20	BU	1159	0	1228	26	0
21	BV	1231	0	1278	30	0
22	BW	1374	0	1405	36	0
23	BX	1181	0	1139	28	0
24	BY	1678	0	1683	32	0
25	Ba	3173	0	3153	68	0
26	Bb	2952	0	2840	61	0
27	Bc	2408	0	2415	61	0
28	Bd	1158	0	1166	37	0
29	Be	972	0	971	25	0
30	Bf	827	0	736	21	0
31	Bg	1167	0	1173	32	0
32	Bh	2319	0	2332	44	0
33	Bi	2138	0	2139	49	0
34	Bj	1775	0	1797	61	0
35	Bk	1246	0	1251	51	0
36	Bl	1097	0	1080	14	0
37	Bm	893	0	878	11	0
38	Bn	837	0	860	28	0
39	Bo	772	0	773	12	0
40	Bp	742	0	749	16	0
41	Bq	474	0	452	16	0
42	Bt	780	0	792	21	0
43	Bu	1198	0	1192	24	0
44	Bv	1131	0	1093	26	0
45	Bw	3126	0	3153	59	0
46	Bx	1325	0	1354	36	0
47	AA	20411	0	10344	285	0
48	CL	317	0	309	11	0
48	DL	213	0	234	8	0
48	EL	222	0	247	6	0
48	FL	213	0	234	11	0
48	GL	213	0	234	7	0
48	HL	205	0	223	8	0
49	AB	1762	0	1774	41	0
50	AC	1075	0	1087	30	0
51	AE	2732	0	2774	52	0
52	AF	981	0	1012	32	0
53	AG	1721	0	1747	47	0
54	AI	2650	0	2613	58	0
55	AJ	1155	0	1193	43	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
56	AK	1007	0	1042	30	0
57	AL	840	0	890	22	0
58	AN	858	0	883	23	0
59	AO	1448	0	1536	24	0
60	AP	932	0	951	35	0
61	AQ	875	0	931	18	0
62	AR	784	0	813	22	0
63	AU	734	0	745	18	0
64	AV	1498	0	766	22	0
65	AX	131	0	66	3	0
66	AZ	90	0	89	0	0
67	Aa	2378	0	2392	69	0
68	Ab	1101	0	1118	17	0
69	Ac	1367	0	1385	27	0
70	Ad	1467	0	1466	30	0
71	Ae	3109	0	3029	94	0
72	Af	778	0	791	19	0
73	Ag	2875	0	2886	75	0
74	Ah	1015	0	969	32	0
75	Ai	824	0	842	27	0
76	Aj	1788	0	1799	67	0
77	Ak	2222	0	2270	72	0
78	Am	930	0	959	20	0
79	An	639	0	709	13	0
80	Ao	4526	0	4434	129	0
81	Ap	1564	0	1531	36	0
82	B0	857	0	878	17	0
83	B1	2036	0	2058	36	0
84	B2	1548	0	1583	33	0
85	B3	968	0	1018	28	0
86	B4	474	0	454	15	0
87	B5	902	0	916	20	0
88	B6	425	0	471	13	0
89	AR	1	0	0	0	0
89	Ac	1	0	0	0	0
89	Ap	1	0	0	0	0
89	B5	1	0	0	0	0
89	B9	1	0	0	0	0
89	Bx	1	0	0	0	0
90	AA	101	0	0	0	0
90	AB	1	0	0	0	0
90	AL	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
90	AV	1	0	0	0	0
90	AX	1	0	0	0	0
90	Ag	1	0	0	0	0
90	Am	1	0	0	0	0
90	An	1	0	0	0	0
90	B3	1	0	0	0	0
90	BA	212	0	0	0	0
90	BB	1	0	0	0	0
90	BD	3	0	0	0	0
90	BE	1	0	0	0	0
90	BP	3	0	0	0	0
90	BQ	1	0	0	0	0
90	BR	1	0	0	0	0
90	Be	2	0	0	0	0
90	Bl	1	0	0	0	0
90	Bt	1	0	0	0	0
91	BA	48	0	23	2	0
92	AA	14	0	26	1	0
92	BA	28	0	52	2	0
93	Ag	32	0	12	3	0
94	Ag	3	0	0	0	0
All	All	174319	0	147376	3062	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 10.

All (3062) 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)
81:Ap:105:CYS:HB2	81:Ap:108:CYS:HB3	1.48	0.95
47:AA:405:G:H1	47:AA:452:C:HO2'	0.90	0.89
76:Aj:53:ARG:HG2	81:Ap:223:ARG:HH12	1.42	0.85
25:Ba:126:THR:HG22	25:Ba:372:ASN:HB2	1.57	0.84
4:BA:293:U:HO2'	4:BA:1434:U:HO2'	1.24	0.84
60:AP:102:MET:HE3	60:AP:106:ASN:HD21	1.42	0.84
4:BA:751:U:OP2	23:BX:50:ARG:NH1	2.11	0.84
80:Ao:265:MET:HB3	80:Ao:269:MET:HE1	1.59	0.84
55:AJ:74:LYS:HD3	55:AJ:75:ARG:HG3	1.59	0.83
74:Ah:284:ARG:NH2	80:Ao:451:ILE:O	2.11	0.83
55:AJ:180:LEU:HD12	55:AJ:181:PRO:HD2	1.62	0.82
18:BS:118:SER:OG	18:BS:121:ASN:OD1	1.96	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BX:51:VAL:HG11	29:Be:50:GLN:HE22	1.45	0.81
26:Bb:335:LEU:HB3	26:Bb:337:MET:HE3	1.61	0.81
22:BW:210:LEU:HD21	31:Bg:121:THR:HG22	1.60	0.81
25:Ba:125:LYS:HG2	25:Ba:371:LYS:HG3	1.61	0.81
80:Ao:420:GLN:HG3	80:Ao:456:ARG:HD2	1.63	0.80
78:Am:76:CYS:HA	78:Am:79:ARG:HE	1.46	0.80
15:BP:119:THR:O	15:BP:123:ASN:ND2	2.15	0.80
50:AC:128:PRO:HD2	50:AC:131:LYS:HD3	1.64	0.79
47:AA:221:A:OP1	81:Ap:148:LYS:NZ	2.14	0.79
53:AG:71:THR:CG2	54:AI:252:GLN:O	2.32	0.78
24:BY:168:GLU:HG3	24:BY:169:THR:HG23	1.65	0.78
60:AP:30:PRO:HB2	60:AP:56:PRO:HG2	1.65	0.78
69:Ac:56:GLN:HG2	69:Ac:64:ILE:HD12	1.64	0.78
16:BQ:218:ILE:HG23	16:BQ:223:MET:HB2	1.66	0.78
10:BJ:237:VAL:HA	48:GL:76:LEU:HD11	1.66	0.77
16:BQ:92:LEU:HD11	16:BQ:188:LYS:HE2	1.65	0.77
71:Ae:70:LEU:HD23	71:Ae:405:LEU:HD11	1.66	0.77
80:Ao:533:GLU:HA	80:Ao:536:MET:HE3	1.65	0.77
28:Bd:72:ILE:HD11	35:Bk:207:LEU:HD22	1.67	0.77
4:BA:472:U:H5''	85:B3:74:SER:HB2	1.65	0.77
56:AK:90:ILE:HG13	56:AK:99:ILE:HG12	1.66	0.76
4:BA:548:A:H5''	41:Bq:126:ILE:HD11	1.65	0.76
6:BD:133:ASP:OD1	6:BD:136:ARG:NH1	2.19	0.76
34:Bj:48:LEU:HD21	34:Bj:175:GLN:HB3	1.66	0.75
4:BA:112:A:N6	4:BA:115:U:OP2	2.19	0.75
34:Bj:152:LYS:HB2	34:Bj:157:LEU:HD21	1.68	0.75
17:BR:130:CYS:HG	87:B5:156:THR:HG1	1.24	0.75
2:B8:107:VAL:HG21	2:B8:161:MET:HE2	1.67	0.75
13:BN:27:PRO:HG2	13:BN:30:LYS:HB2	1.68	0.75
47:AA:50:U:H2'	47:AA:51:G:H8	1.51	0.75
4:BA:11:G:H5'	84:B2:228:LEU:HD11	1.67	0.74
45:Bw:146:GLN:HA	45:Bw:150:PHE:HB2	1.69	0.74
60:AP:110:LEU:HA	60:AP:113:LYS:HD3	1.68	0.74
67:Aa:100:ILE:HG23	67:Aa:321:SER:HB2	1.68	0.74
74:Ah:284:ARG:HD3	74:Ah:285:PRO:HD2	1.70	0.74
10:BJ:64:CYS:SG	46:Bx:70:CYS:HB2	2.27	0.74
23:BX:10:TYR:HA	23:BX:14:ASN:HD22	1.52	0.74
7:BE:107:MET:HE3	19:BT:95:GLU:HG3	1.69	0.73
24:BY:78:GLN:OE1	33:Bi:74:ARG:NH1	2.21	0.73
26:Bb:241:PRO:HB2	26:Bb:244:ARG:HH21	1.53	0.73
4:BA:529:G:H1'	4:BA:549:C:H1'	1.70	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:BA:1569:A:N1	7:BE:69:ASN:ND2	2.36	0.73
8:BF:262:THR:HG22	8:BF:264:PRO:HD2	1.69	0.73
47:AA:16:U:O2'	47:AA:294:A:N6	2.22	0.73
16:BQ:33:LEU:HD21	41:Bq:122:ARG:HG3	1.71	0.73
85:B3:69:VAL:HG22	85:B3:119:ILE:HG12	1.71	0.73
5:BB:35:G:N2	35:Bk:124:SER:O	2.22	0.72
51:AE:304:LYS:HE2	51:AE:409:GLU:HG2	1.70	0.72
71:Ae:119:ILE:HB	71:Ae:155:GLN:HE21	1.54	0.72
12:BL:97:SER:O	12:BL:146:ASN:ND2	2.22	0.72
60:AP:51:LEU:HD13	60:AP:73:ILE:HG12	1.72	0.72
71:Ae:110:LEU:HD11	71:Ae:431:ILE:HD13	1.71	0.72
80:Ao:543:HIS:HB2	80:Ao:588:ARG:HD2	1.71	0.72
23:BX:37:GLU:HB3	23:BX:104:THR:HG23	1.71	0.72
71:Ae:157:LEU:HA	71:Ae:161:ALA:HB2	1.71	0.72
4:BA:834:C:H41	33:Bi:39:LYS:HE2	1.55	0.72
60:AP:53:SER:H	60:AP:67:ALA:HB3	1.55	0.72
59:AO:110:MET:SD	59:AO:114:ASN:ND2	2.63	0.71
69:Ac:51:ASN:ND2	69:Ac:95:ASN:OD1	2.23	0.71
22:BW:151:LEU:HB2	22:BW:167:LYS:HB2	1.71	0.71
67:Aa:205:LYS:NZ	81:Ap:182:GLY:O	2.23	0.71
4:BA:706:A:OP1	23:BX:50:ARG:NH2	2.22	0.71
10:BJ:95:MET:HB2	10:BJ:159:PRO:HB3	1.72	0.71
67:Aa:287:THR:HG23	67:Aa:289:HIS:H	1.55	0.71
87:B5:138:ARG:HA	87:B5:141:MET:HG2	1.70	0.71
34:Bj:140:ALA:HA	34:Bj:143:LYS:HE3	1.71	0.71
25:Ba:283:TYR:OH	45:Bw:146:GLN:NE2	2.23	0.71
56:AK:130:ALA:HB1	56:AK:163:ALA:HB2	1.73	0.71
18:BS:122:VAL:HG13	18:BS:157:SER:HA	1.71	0.71
4:BA:1492:C:H5''	19:BT:271:ARG:HD3	1.72	0.71
24:BY:146:VAL:HG12	24:BY:153:ILE:HG22	1.73	0.71
51:AE:305:MET:HE3	51:AE:332:LEU:HD21	1.73	0.71
43:Bu:40:GLN:HB3	43:Bu:47:LYS:HZ1	1.56	0.70
46:Bx:72:ILE:HD11	46:Bx:115:ILE:HD11	1.71	0.70
67:Aa:211:ARG:O	67:Aa:215:ILE:HG12	1.91	0.70
74:Ah:284:ARG:NH2	80:Ao:448:TRP:O	2.24	0.70
36:Bl:96:VAL:HG22	36:Bl:110:ILE:HG12	1.74	0.70
47:AA:881:A:H5''	71:Ae:438:ARG:HH12	1.57	0.70
73:Ag:122:ARG:NH2	73:Ag:296:VAL:O	2.24	0.70
35:Bk:175:HIS:HA	35:Bk:178:LEU:HG	1.72	0.70
34:Bj:256:SER:O	34:Bj:260:LEU:HB2	1.92	0.70
4:BA:392:U:H2'	4:BA:393:A:H8	1.56	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
47:AA:227:A:H61	47:AA:270:C:H42	1.40	0.70
73:Ag:59:ARG:HB2	73:Ag:62:HIS:HD2	1.56	0.70
26:Bb:124:ARG:NH2	28:Bd:112:GLU:OE1	2.25	0.69
51:AE:150:ARG:HB3	51:AE:155:GLN:HE21	1.57	0.69
71:Ae:169:LEU:HG	71:Ae:206:GLU:HG3	1.73	0.69
77:Ak:171:ARG:NH2	77:Ak:212:ASP:OD1	2.24	0.69
76:Aj:70:ARG:HH21	76:Aj:141:LEU:HD21	1.56	0.69
80:Ao:419:PHE:HE2	80:Ao:450:LEU:HD13	1.58	0.69
14:BO:104:ASN:HD21	19:BT:158:GLN:HE21	1.41	0.69
91:BA:3212:5GP:OP1	21:BV:182:LYS:NZ	2.25	0.69
50:AC:163:VAL:HG11	50:AC:167:ILE:HB	1.74	0.69
71:Ae:370:GLN:HB3	71:Ae:374:ARG:HH12	1.57	0.69
4:BA:550:A:N7	4:BA:551:A:N6	2.40	0.69
35:Bk:94:HIS:ND1	35:Bk:154:GLU:OE2	2.21	0.69
35:Bk:94:HIS:HB2	35:Bk:185:SER:HB3	1.75	0.69
77:Ak:63:ALA:HB3	77:Ak:66:GLN:HB2	1.75	0.69
22:BW:73:HIS:HD2	22:BW:118:VAL:HG13	1.57	0.68
34:Bj:264:LEU:HD13	34:Bj:269:LEU:HB3	1.75	0.68
26:Bb:221:LEU:HB2	26:Bb:267:ARG:HG3	1.74	0.68
47:AA:58:U:N3	76:Aj:57:ASP:OD1	2.17	0.68
8:BF:132:GLY:HA3	33:Bi:42:THR:HG21	1.75	0.68
67:Aa:240:MET:HE3	67:Aa:249:VAL:HG22	1.75	0.68
27:Bc:175:ALA:HB2	27:Bc:316:SER:HB2	1.76	0.68
4:BA:803:A:N3	4:BA:805:C:N4	2.40	0.68
15:BP:225:ASP:HB3	15:BP:228:ARG:HG2	1.75	0.68
27:Bc:78:MET:HE3	27:Bc:83:SER:HB3	1.76	0.68
22:BW:206:ILE:HG22	22:BW:209:ALA:HB2	1.76	0.68
53:AG:71:THR:HG21	54:AI:252:GLN:HB3	1.74	0.68
71:Ae:358:LEU:HB2	71:Ae:361:GLU:HG3	1.76	0.68
8:BF:49:ARG:NH2	8:BF:270:GLU:OE1	2.25	0.68
10:BJ:46:LYS:HE3	42:Bt:31:ALA:HB1	1.76	0.68
29:Be:121:GLN:HE22	29:Be:128:LEU:HD12	1.59	0.68
45:Bw:236:ARG:HA	45:Bw:289:GLY:HA2	1.76	0.68
72:Af:105:ILE:HG12	72:Af:115:ILE:HG12	1.74	0.68
80:Ao:67:LYS:HG3	80:Ao:68:VAL:HG13	1.76	0.68
80:Ao:249:ARG:HA	80:Ao:253:LEU:HD23	1.76	0.68
24:BY:133:VAL:HG11	24:BY:145:ARG:HD2	1.74	0.68
14:BO:133:GLU:HG3	14:BO:134:PHE:HD1	1.59	0.67
47:AA:386:U:H4'	52:AF:94:VAL:HG12	1.76	0.67
76:Aj:41:MET:HE1	76:Aj:55:TRP:HA	1.76	0.67
49:AB:101:MET:HE2	49:AB:228:PRO:HD3	1.76	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
11:BK:35:LEU:HD12	11:BK:38:ILE:HD11	1.76	0.67
15:BP:10:PRO:HB2	15:BP:13:LEU:HD23	1.74	0.67
32:Bh:97:TYR:HB2	32:Bh:181:SER:HA	1.75	0.67
47:AA:492:U:OP1	79:An:139:ASN:ND2	2.28	0.67
49:AB:215:ASN:OD1	68:Ab:15:ARG:NH1	2.28	0.67
34:Bj:85:SER:HB2	86:B4:51:ARG:HB3	1.76	0.67
53:AG:119:LYS:NZ	73:Ag:364:TRP:O	2.26	0.67
71:Ae:320:LEU:HD13	71:Ae:323:LEU:HD12	1.76	0.67
4:BA:831:U:OP2	4:BA:836:A:N6	2.24	0.67
71:Ae:96:LEU:HA	71:Ae:99:LYS:HZ3	1.59	0.67
4:BA:61:U:OP2	83:B1:100:ARG:NH2	2.25	0.67
33:Bi:267:PRO:HG2	33:Bi:270:ALA:HB2	1.76	0.67
10:BJ:47:LEU:HD22	16:BQ:226:ILE:HG12	1.77	0.67
74:Ah:350:LEU:O	74:Ah:361:LYS:NZ	2.24	0.67
25:Ba:160:HIS:HA	25:Ba:164:TRP:HB2	1.76	0.67
35:Bk:137:LEU:HD21	35:Bk:141:GLY:HA2	1.75	0.67
47:AA:29:A:H2'	47:AA:30:A:H8	1.60	0.67
59:AO:144:MET:HE1	59:AO:150:ASP:HB3	1.77	0.67
22:BW:82:LYS:HD3	22:BW:147:ARG:HB3	1.77	0.66
51:AE:92:LYS:HE3	75:Ai:75:HIS:HB3	1.76	0.66
74:Ah:332:ASP:HB3	74:Ah:369:ARG:HH12	1.60	0.66
45:Bw:195:ASP:HB2	45:Bw:234:GLN:HB3	1.78	0.66
10:BJ:154:LEU:HD13	10:BJ:164:MET:HE2	1.76	0.66
47:AA:11:G:H1	47:AA:524:U:H3	1.42	0.66
24:BY:161:ARG:HH21	24:BY:165:ILE:HB	1.61	0.66
24:BY:191:LEU:HG	29:Be:93:THR:HG21	1.78	0.66
4:BA:789:A:N6	4:BA:1000:A:O2'	2.28	0.66
7:BE:45:SER:O	27:Bc:303:ARG:NH2	2.28	0.66
29:Be:16:ASP:OD1	29:Be:25:ARG:NH2	2.28	0.66
71:Ae:121:SER:HA	71:Ae:159:TYR:HE2	1.58	0.66
4:BA:849:U:P	6:BD:288:ARG:HH22	2.19	0.66
50:AC:165:LYS:HG3	50:AC:166:TYR:H	1.59	0.66
74:Ah:379:LEU:HD13	74:Ah:386:LEU:HB2	1.76	0.66
80:Ao:353:LYS:HB3	80:Ao:381:LEU:HD22	1.76	0.66
26:Bb:50:LYS:HG3	82:B0:121:PRO:HA	1.77	0.66
7:BE:104:LEU:HB2	7:BE:121:LEU:HB2	1.77	0.66
9:BI:120:ARG:NH1	83:B1:132:LEU:HB3	2.11	0.66
51:AE:191:ARG:NH2	81:Ap:80:ASN:OD1	2.29	0.66
4:BA:1026:A:O2'	4:BA:1277:G:N2	2.23	0.65
8:BF:60:ARG:HG2	8:BF:61:PRO:HD2	1.77	0.65
16:BQ:49:LYS:HA	16:BQ:113:MET:HE1	1.77	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:Ba:360:ASN:HD22	25:Ba:372:ASN:HA	1.61	0.65
60:AP:35:VAL:HG23	60:AP:50:GLN:HA	1.77	0.65
4:BA:95:U:OP1	44:Bv:63:ARG:NH2	2.29	0.65
4:BA:1207:C:O2'	82:B0:88:CYS:SG	2.54	0.65
47:AA:202:A:N1	76:Aj:53:ARG:NH2	2.43	0.65
47:AA:883:C:H3'	71:Ae:433:ARG:HH12	1.62	0.65
53:AG:71:THR:HG22	54:AI:252:GLN:O	1.96	0.65
4:BA:198:U:H2'	4:BA:199:A:H8	1.62	0.65
15:BP:19:LEU:HD13	42:Bt:95:LEU:HD12	1.78	0.65
20:BU:33:GLY:O	20:BU:36:ASN:ND2	2.30	0.65
57:AL:62:VAL:HA	57:AL:83:VAL:HG12	1.78	0.65
4:BA:569:C:N4	4:BA:1020:C:O2	2.30	0.65
18:BS:138:GLU:OE2	26:Bb:136:ARG:NH1	2.30	0.65
34:Bj:196:SER:O	34:Bj:198:ASN:ND2	2.29	0.65
83:B1:172:GLN:HA	83:B1:188:TYR:HE2	1.61	0.65
5:BB:32:A:OP1	86:B4:43:ARG:NH2	2.30	0.65
27:Bc:112:MET:HB2	27:Bc:267:PRO:HB3	1.78	0.65
44:Bv:40:PRO:HG3	44:Bv:51:GLN:HB3	1.78	0.65
8:BF:59:ARG:HE	8:BF:84:PRO:HB2	1.62	0.65
38:Bn:35:ARG:NH1	38:Bn:37:THR:O	2.30	0.65
33:Bi:202:LEU:HD22	33:Bi:205:GLN:HG2	1.79	0.65
80:Ao:269:MET:HA	80:Ao:272:HIS:HB2	1.77	0.65
77:Ak:154:ASP:HB2	77:Ak:174:THR:HB	1.77	0.65
16:BQ:204:GLU:OE1	16:BQ:207:ARG:NH2	2.30	0.65
51:AE:203:LEU:HD22	51:AE:222:ILE:HD11	1.79	0.65
24:BY:197:GLU:HG2	84:B2:97:ASN:HB2	1.77	0.64
34:Bj:235:LYS:NZ	35:Bk:172:GLU:OE2	2.31	0.64
49:AB:106:PHE:HB2	49:AB:116:ASP:HB3	1.79	0.64
76:Aj:53:ARG:HG2	81:Ap:223:ARG:NH1	2.10	0.64
49:AB:238:ASN:ND2	49:AB:241:SER:OG	2.30	0.64
52:AF:49:TYR:OH	52:AF:90:ARG:NH1	2.30	0.64
27:Bc:186:LEU:O	27:Bc:291:ARG:NH2	2.29	0.64
35:Bk:178:LEU:HD21	35:Bk:184:LEU:HD12	1.78	0.64
46:Bx:39:VAL:HG11	46:Bx:52:ARG:HH22	1.63	0.64
4:BA:129:A:OP1	8:BF:142:ARG:NH2	2.30	0.64
7:BE:134:PRO:HG2	7:BE:137:ASN:HB2	1.79	0.64
7:BE:248:VAL:HG12	7:BE:250:ARG:HG2	1.80	0.64
13:BN:152:PRO:HB3	46:Bx:84:GLU:HG3	1.80	0.64
20:BU:78:GLU:OE2	31:Bg:135:ASN:ND2	2.30	0.64
47:AA:344:G:O6	56:AK:124:LYS:NZ	2.29	0.64
4:BA:975:G:O2'	4:BA:977:G:OP2	2.15	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:BA:1014:A:OP1	20:BU:34:ARG:NH2	2.30	0.64
8:BF:66:ILE:HD13	8:BF:263:LEU:HD11	1.78	0.64
25:Ba:135:LYS:HA	25:Ba:138:ARG:HE	1.63	0.64
26:Bb:307:HIS:HB2	26:Bb:311:MET:HE2	1.78	0.64
4:BA:602:G:N7	15:BP:44:ARG:NH2	2.43	0.64
4:BA:737:G:O4'	25:Ba:302:ARG:NH1	2.31	0.64
4:BA:743:U:H3'	4:BA:744:A:H5''	1.78	0.64
6:BD:178:ARG:HG2	6:BD:179:GLU:H	1.63	0.64
28:Bd:67:ARG:HB3	28:Bd:71:ARG:HH12	1.62	0.64
34:Bj:215:PHE:HB3	34:Bj:268:TYR:CE2	2.33	0.64
81:Ap:68:TYR:OH	81:Ap:110:ASP:OD2	2.16	0.64
4:BA:106:C:OP1	32:Bh:33:LYS:NZ	2.31	0.64
7:BE:145:LEU:HD12	7:BE:181:VAL:HG21	1.80	0.64
7:BE:275:ARG:HD2	7:BE:333:TYR:HE1	1.63	0.64
9:BI:105:LEU:HD12	9:BI:112:VAL:HG11	1.79	0.64
4:BA:486:A:H61	4:BA:582:C:H42	1.46	0.64
4:BA:148:A:OP1	32:Bh:52:ARG:NH1	2.31	0.64
13:BN:13:THR:HG21	22:BW:206:ILE:HD11	1.80	0.64
2:B8:113:ARG:NH2	4:BA:84:G:OP2	2.31	0.64
16:BQ:163:MET:HE1	16:BQ:175:PHE:HZ	1.63	0.64
19:BT:96:ARG:NH1	19:BT:285:GLU:OE1	2.31	0.64
26:Bb:45:LEU:O	26:Bb:50:LYS:NZ	2.31	0.64
39:Bo:74:ALA:HB2	85:B3:137:THR:HG21	1.79	0.64
4:BA:1168:A:N1	26:Bb:27:ARG:N	2.46	0.63
8:BF:79:LEU:HD23	37:Bm:95:VAL:HG13	1.80	0.63
24:BY:25:ARG:HB3	24:BY:28:SER:HB3	1.78	0.63
53:AG:71:THR:HG21	54:AI:252:GLN:O	1.98	0.63
27:Bc:169:VAL:HG12	27:Bc:322:GLU:HB3	1.80	0.63
27:Bc:224:LEU:HD13	27:Bc:234:LEU:HD22	1.79	0.63
31:Bg:65:VAL:HB	37:Bm:155:ILE:HG23	1.80	0.63
84:B2:71:ASP:OD1	84:B2:72:ASP:N	2.29	0.63
4:BA:470:C:OP2	85:B3:77:ARG:NH1	2.32	0.63
28:Bd:87:ASP:OD1	35:Bk:153:HIS:CE1	2.51	0.63
47:AA:800:G:H2'	47:AA:801:A:C8	2.33	0.63
50:AC:136:VAL:HG23	50:AC:153:LEU:HD23	1.81	0.63
7:BE:210:THR:OG1	7:BE:262:GLY:O	2.15	0.63
43:Bu:80:TYR:HE1	43:Bu:146:LEU:HD21	1.63	0.63
54:AI:115:GLY:HA3	55:AJ:84:ASP:HB2	1.80	0.63
64:AV:4:A:H2'	64:AV:5:A:C8	2.33	0.63
4:BA:869:G:O2'	4:BA:966:U:OP2	2.17	0.63
30:Bf:68:PRO:HG2	30:Bf:71:HIS:HD2	1.63	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
54:AI:311:ARG:HA	54:AI:314:LEU:HD12	1.79	0.63
60:AP:20:ARG:NH2	60:AP:41:CYS:SG	2.71	0.63
60:AP:67:ALA:HB2	67:Aa:218:TYR:CZ	2.33	0.63
4:BA:528:U:O3'	41:Bq:115:ARG:NH2	2.30	0.63
4:BA:877:U:H5''	4:BA:878:G:H5'	1.80	0.63
7:BE:236:THR:HA	7:BE:239:ARG:HG3	1.80	0.63
15:BP:142:GLU:HG3	15:BP:162:LEU:HD23	1.80	0.63
31:Bg:44:ARG:O	31:Bg:48:GLU:HG3	1.98	0.63
67:Aa:244:ASP:HB2	67:Aa:283:LEU:HD11	1.79	0.63
35:Bk:189:HIS:HE1	35:Bk:194:PHE:HB2	1.64	0.63
42:Bt:45:ASN:HD21	85:B3:118:ARG:HA	1.64	0.63
55:AJ:121:LEU:HB2	58:AN:108:ARG:HE	1.64	0.63
69:Ac:78:PHE:HB2	69:Ac:86:VAL:HB	1.81	0.63
77:Ak:213:ARG:NH1	77:Ak:223:TYR:OH	2.32	0.63
6:BD:203:ARG:HH21	6:BD:206:GLN:HE22	1.46	0.63
6:BD:270:ARG:HG3	6:BD:271:PRO:HD2	1.80	0.63
18:BS:75:ARG:NH2	18:BS:97:ARG:O	2.31	0.63
21:BV:198:ARG:NH2	31:Bg:13:SER:OG	2.32	0.63
52:AF:27:GLU:HG3	70:Ad:170:ARG:HH12	1.64	0.63
54:AI:397:ARG:NH2	64:AV:30:C:OP2	2.31	0.63
61:AQ:95:VAL:HG23	61:AQ:96:THR:HG23	1.79	0.63
10:BJ:97:ALA:HB3	10:BJ:154:LEU:HB2	1.80	0.63
34:Bj:47:LEU:HD13	34:Bj:178:TRP:CE2	2.34	0.63
36:Bl:138:THR:HG22	36:Bl:149:VAL:HG12	1.80	0.63
40:Bp:76:MET:HG2	46:Bx:49:ILE:HB	1.80	0.63
45:Bw:53:ALA:O	45:Bw:62:ARG:NH2	2.31	0.63
58:AN:50:GLU:O	58:AN:54:ILE:HG12	1.98	0.63
65:AX:2:A:O2'	65:AX:3:A:OP1	2.16	0.63
71:Ae:136:ARG:NH1	71:Ae:173:VAL:O	2.32	0.63
6:BD:76:PRO:HB2	6:BD:103:ARG:HD2	1.79	0.62
20:BU:54:THR:HG21	21:BV:176:MET:H	1.64	0.62
25:Ba:35:VAL:HB	25:Ba:38:THR:HG22	1.80	0.62
69:Ac:90:VAL:HG23	69:Ac:93:LYS:HB2	1.81	0.62
78:Am:100:PRO:HD2	78:Am:103:LYS:HD3	1.81	0.62
86:B4:60:GLN:NE2	86:B4:79:PRO:O	2.32	0.62
7:BE:69:ASN:HD21	7:BE:154:HIS:HE1	1.46	0.62
29:Be:88:GLN:O	29:Be:92:GLU:HG2	2.00	0.62
47:AA:90:U:O5'	70:Ad:68:ARG:NH2	2.32	0.62
80:Ao:531:LYS:HE2	80:Ao:562:TYR:HE2	1.64	0.62
54:AI:136:ARG:HA	54:AI:136:ARG:HH11	1.64	0.62
4:BA:875:C:O2'	79:An:151:ARG:NH2	2.31	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:BE:275:ARG:HD2	7:BE:333:TYR:CE1	2.34	0.62
45:Bw:147:GLU:HG3	45:Bw:170:PHE:HA	1.80	0.62
47:AA:957:A:H5''	78:Am:24:ASN:HD21	1.64	0.62
67:Aa:307:LEU:HD21	67:Aa:350:ILE:HG23	1.81	0.62
70:Ad:112:GLU:OE2	70:Ad:115:ARG:NH2	2.32	0.62
1:B7:58:PRO:HB2	4:BA:672:G:H4'	1.80	0.62
4:BA:703:A:OP1	23:BX:41:GLN:NE2	2.32	0.62
6:BD:162:ASP:OD1	6:BD:163:THR:N	2.31	0.62
47:AA:66:C:H42	70:Ad:27:ARG:HB2	1.65	0.62
4:BA:1354:A:H61	4:BA:1465:A:H61	1.46	0.62
16:BQ:109:ILE:O	16:BQ:113:MET:HB3	1.99	0.62
10:BJ:44:ARG:NH2	16:BQ:234:ASP:OD1	2.30	0.62
4:BA:35:A:OP1	29:Be:32:LYS:NZ	2.30	0.62
4:BA:1454:C:H2'	4:BA:1455:C:C6	2.34	0.62
47:AA:876:A:H5'	47:AA:877:A:H5'	1.80	0.62
59:AO:176:TYR:HB2	61:AQ:89:GLY:HA3	1.80	0.62
4:BA:1521:U:H5	13:BN:177:ARG:HB3	1.64	0.62
8:BF:84:PRO:HA	8:BF:88:SER:HB3	1.80	0.62
47:AA:693:A:OP2	75:Ai:89:ARG:NH2	2.33	0.62
64:AV:40:A:H2'	64:AV:41:A:C8	2.35	0.62
4:BA:266:C:O2'	4:BA:281:A:N6	2.33	0.62
8:BF:63:GLN:NE2	8:BF:81:GLU:OE2	2.33	0.62
12:BL:178:ILE:HG22	12:BL:182:ASN:HD21	1.63	0.62
26:Bb:279:ILE:HD11	26:Bb:310:ALA:HB3	1.82	0.62
32:Bh:60:ARG:HD2	32:Bh:63:LYS:HD2	1.82	0.62
80:Ao:586:PHE:HD2	80:Ao:591:ARG:HB3	1.65	0.62
22:BW:105:GLU:OE1	22:BW:116:LYS:NZ	2.28	0.61
54:AI:148:HIS:HE1	54:AI:160:SER:HB3	1.65	0.61
11:BK:122:ARG:O	11:BK:126:GLN:NE2	2.30	0.61
13:BN:117:HIS:O	13:BN:121:MET:HG2	2.00	0.61
15:BP:137:GLY:HA3	15:BP:157:GLN:HG3	1.82	0.61
31:Bg:133:PHE:HZ	32:Bh:261:SER:HB2	1.65	0.61
35:Bk:170:PHE:O	35:Bk:174:ILE:HG12	2.00	0.61
73:Ag:105:LEU:HD23	73:Ag:108:LEU:HD21	1.82	0.61
73:Ag:259:VAL:HB	73:Ag:305:ILE:HD13	1.81	0.61
85:B3:104:PRO:HA	85:B3:107:ASN:HB2	1.82	0.61
4:BA:1404:G:N2	4:BA:1404:G:OP2	2.32	0.61
33:Bi:154:ASN:HD22	33:Bi:154:ASN:C	2.06	0.61
68:Ab:89:ASN:HB3	68:Ab:92:PHE:HB2	1.80	0.61
74:Ah:326:LEU:HB2	74:Ah:347:THR:HG23	1.82	0.61
4:BA:1570:C:H3'	4:BA:1571:A:H2'	1.81	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:BB:63:C:H2'	5:BB:64:A:H8	1.65	0.61
43:Bu:96:SER:O	43:Bu:145:ASN:ND2	2.33	0.61
46:Bx:99:MET:HE3	46:Bx:100:LEU:H	1.65	0.61
47:AA:918:A:H2'	47:AA:919:A:H5''	1.82	0.61
51:AE:93:LEU:H	75:Ai:75:HIS:HD2	1.47	0.61
70:Ad:62:GLN:OE1	70:Ad:65:GLN:NE2	2.33	0.61
80:Ao:137:ILE:HD12	80:Ao:138:PRO:HD2	1.82	0.61
38:Bn:77:PRO:HG2	38:Bn:80:LEU:HB2	1.81	0.61
47:AA:283:U:OP1	51:AE:419:ARG:NH1	2.33	0.61
60:AP:59:ASN:HD21	60:AP:63:GLU:HG2	1.64	0.61
62:AR:83:GLN:HE22	70:Ad:191:ILE:HG12	1.65	0.61
71:Ae:270:LEU:HD12	71:Ae:273:LYS:HD3	1.81	0.61
4:BA:83:A:OP2	4:BA:1235:C:O2'	2.16	0.61
4:BA:1278:C:H2'	4:BA:1279:G:H8	1.65	0.61
55:AJ:123:SER:HB3	55:AJ:127:PHE:HB2	1.82	0.61
69:Ac:140:MET:HE2	69:Ac:146:GLN:HG2	1.82	0.61
4:BA:22:U:O4	84:B2:172:ARG:NH1	2.33	0.61
64:AV:48:G:H2'	64:AV:49:G:C8	2.36	0.61
73:Ag:99:MET:H	93:Ag:500:GTP:HN1	1.49	0.61
80:Ao:407:ARG:HG3	80:Ao:445:GLY:HA2	1.83	0.61
4:BA:1158:C:O2'	4:BA:1251:C:OP2	2.17	0.61
10:BJ:221:MET:HE1	48:FL:83:ASN:HA	1.82	0.61
28:Bd:150:LEU:HB2	34:Bj:212:HIS:CE1	2.35	0.61
60:AP:29:ARG:NH2	60:AP:55:ASP:OD1	2.31	0.61
64:AV:4:A:H2'	64:AV:5:A:H8	1.66	0.61
67:Aa:128:MET:HB3	67:Aa:132:GLN:HB2	1.83	0.61
67:Aa:250:LEU:HD11	67:Aa:271:THR:HG21	1.81	0.61
73:Ag:59:ARG:HB2	73:Ag:62:HIS:CD2	2.35	0.61
41:Bq:106:LYS:HA	41:Bq:106:LYS:HE3	1.82	0.61
59:AO:102:LYS:O	59:AO:106:LYS:HG3	2.01	0.61
4:BA:971:A:H5''	79:An:163:ARG:HH12	1.66	0.60
8:BF:160:SER:HB3	38:Bn:80:LEU:HD13	1.83	0.60
28:Bd:171:PHE:HB3	35:Bk:186:VAL:HB	1.83	0.60
55:AJ:158:GLU:HG2	80:Ao:70:ILE:HD13	1.83	0.60
73:Ag:323:LEU:HD13	77:Ak:306:ASP:HB2	1.83	0.60
34:Bj:159:LEU:HD11	34:Bj:252:HIS:HB2	1.82	0.60
47:AA:162:C:OP2	47:AA:163:A:N6	2.34	0.60
47:AA:880:U:H2'	47:AA:881:A:C8	2.36	0.60
77:Ak:144:LYS:O	77:Ak:148:HIS:ND1	2.23	0.60
4:BA:243:A:N3	4:BA:1266:U:O2'	2.34	0.60
8:BF:56:PRO:HG2	8:BF:59:ARG:HB2	1.83	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
51:AE:283:GLU:O	51:AE:356:GLN:NE2	2.34	0.60
69:Ac:29:VAL:HB	69:Ac:79:TYR:HB2	1.84	0.60
73:Ag:221:LEU:HD21	73:Ag:242:VAL:HG13	1.82	0.60
4:BA:164:A:OP1	20:BU:52:LYS:NZ	2.33	0.60
7:BE:308:LYS:NZ	7:BE:309:ASP:OD1	2.33	0.60
14:BO:85:LEU:HD23	14:BO:137:VAL:HG11	1.84	0.60
20:BU:85:ALA:O	20:BU:89:ASN:ND2	2.34	0.60
35:Bk:90:VAL:HG12	35:Bk:160:SER:HA	1.83	0.60
35:Bk:92:ASN:OD1	35:Bk:158:GLN:NE2	2.35	0.60
60:AP:42:PRO:HG2	60:AP:45:GLY:HA3	1.83	0.60
70:Ad:115:ARG:O	70:Ad:119:LEU:HD12	2.00	0.60
72:Af:141:VAL:HA	72:Af:173:LEU:HA	1.83	0.60
4:BA:1195:A:OP2	26:Bb:364:ARG:NE	2.31	0.60
4:BA:1226:C:OP1	4:BA:1227:U:N3	2.34	0.60
11:BK:17:SER:N	11:BK:72:VAL:O	2.34	0.60
47:AA:164:C:OP1	76:Aj:19:ARG:NH2	2.32	0.60
77:Ak:92:PRO:HB3	77:Ak:98:PRO:HD3	1.84	0.60
78:Am:101:PRO:O	78:Am:105:ASN:ND2	2.35	0.60
8:BF:212:TRP:O	8:BF:258:THR:OG1	2.19	0.60
10:BJ:82:LEU:HD13	10:BJ:130:VAL:HG21	1.83	0.60
19:BT:152:ARG:HH22	19:BT:191:ARG:HA	1.66	0.60
27:Bc:163:LEU:O	27:Bc:230:ASN:ND2	2.33	0.60
28:Bd:87:ASP:HB3	86:B4:46:ARG:HH22	1.66	0.60
47:AA:958:U:OP2	78:Am:24:ASN:ND2	2.35	0.60
49:AB:140:ILE:HG13	49:AB:187:ARG:HH21	1.65	0.60
62:AR:66:CYS:HB3	62:AR:101:CYS:H	1.65	0.60
4:BA:229:A:H4'	4:BA:230:G:H5'	1.83	0.60
4:BA:376:A:N6	4:BA:1061:U:O2	2.35	0.60
6:BD:258:ILE:HG23	6:BD:263:ARG:HB3	1.83	0.60
12:BL:115:VAL:O	12:BL:119:ILE:HG13	2.02	0.60
15:BP:233:ARG:HH21	15:BP:244:LEU:HD11	1.67	0.60
23:BX:2:ALA:N	23:BX:7:TYR:HH	1.99	0.60
75:Ai:40:GLN:NE2	75:Ai:44:LYS:O	2.34	0.60
4:BA:21:C:H41	83:B1:6:VAL:HG12	1.66	0.60
13:BN:69:GLY:HA2	46:Bx:152:THR:HA	1.83	0.60
14:BO:133:GLU:HG3	14:BO:134:PHE:CD1	2.37	0.60
47:AA:57:A:C8	76:Aj:51:PRO:HB3	2.36	0.60
47:AA:248:C:N4	57:AL:117:ASP:OD2	2.34	0.60
4:BA:407:A:OP2	35:Bk:49:ARG:NH1	2.33	0.60
42:Bt:51:ARG:NH2	85:B3:59:ASP:O	2.34	0.60
47:AA:463:U:O4	47:AA:464:A:N6	2.35	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
49:AB:150:PHE:CD2	49:AB:245:VAL:HG21	2.37	0.60
53:AG:112:THR:HG23	73:Ag:396:TYR:CD1	2.37	0.60
78:Am:10:LEU:O	78:Am:15:ASN:ND2	2.34	0.60
2:B8:110:VAL:O	2:B8:114:PHE:HB2	2.02	0.60
22:BW:136:SER:OG	27:Bc:90:MET:O	2.19	0.60
47:AA:849:G:H5'	79:An:138:MET:HE3	1.84	0.60
71:Ae:173:VAL:HG23	71:Ae:174:GLN:HG3	1.83	0.60
75:Ai:81:GLU:OE1	75:Ai:82:GLN:NE2	2.35	0.60
12:BL:176:HIS:O	12:BL:180:ILE:HG12	2.01	0.59
21:BV:71:PRO:HB2	21:BV:76:GLU:HG2	1.84	0.59
49:AB:197:THR:HG1	49:AB:226:CYS:HG	1.49	0.59
51:AE:374:ARG:HD3	67:Aa:123:PRO:HB3	1.84	0.59
54:AI:365:MET:HG2	54:AI:370:LEU:HB2	1.83	0.59
67:Aa:258:GLU:OE1	67:Aa:259:PRO:HD2	2.02	0.59
4:BA:642:G:OP1	87:B5:93:ARG:NH2	2.35	0.59
4:BA:735:A:OP1	6:BD:103:ARG:NE	2.34	0.59
4:BA:1369:U:H5''	4:BA:1370:U:H5'	1.84	0.59
6:BD:125:GLU:OE1	6:BD:150:ARG:NH2	2.35	0.59
19:BT:240:ILE:HG21	19:BT:243:ILE:HD12	1.84	0.59
52:AF:42:LEU:HB3	52:AF:45:ARG:HH21	1.65	0.59
76:Aj:61:GLU:HG3	76:Aj:83:LYS:HD2	1.83	0.59
77:Ak:254:ILE:HG22	77:Ak:256:GLU:H	1.67	0.59
80:Ao:491:PRO:HB3	80:Ao:495:THR:HB	1.83	0.59
80:Ao:608:ILE:HG21	80:Ao:642:LEU:HD21	1.83	0.59
4:BA:582:C:OP1	20:BU:99:ARG:NH2	2.35	0.59
4:BA:1292:A:N6	4:BA:1304:G:O2'	2.36	0.59
7:BE:275:ARG:HB3	7:BE:284:TYR:CD2	2.38	0.59
75:Ai:66:ARG:NH2	75:Ai:77:ASP:OD1	2.33	0.59
78:Am:88:SER:O	78:Am:92:ASN:ND2	2.35	0.59
85:B3:72:ILE:HD11	85:B3:118:ARG:HB2	1.84	0.59
9:BI:120:ARG:NH1	83:B1:136:ASP:OD1	2.36	0.59
10:BJ:214:LEU:HD13	48:DL:90:LEU:HD23	1.82	0.59
59:AO:181:LYS:O	59:AO:185:GLU:HG2	2.01	0.59
70:Ad:45:PRO:HD2	76:Aj:52:MET:HE2	1.83	0.59
71:Ae:80:TRP:HH2	71:Ae:151:THR:HG23	1.66	0.59
23:BX:30:ARG:NH1	84:B2:120:GLU:OE2	2.34	0.59
39:Bo:45:GLU:OE2	39:Bo:70:ARG:NH2	2.36	0.59
4:BA:851:G:N7	6:BD:231:SER:OG	2.35	0.59
27:Bc:44:LEU:HD21	27:Bc:225:GLU:HG3	1.84	0.59
31:Bg:78:GLU:HG2	31:Bg:84:VAL:HG22	1.85	0.59
46:Bx:72:ILE:HG23	46:Bx:77:LEU:HB2	1.85	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
67:Aa:104:MET:HE3	67:Aa:290:PHE:HE2	1.68	0.59
4:BA:130:A:OP1	8:BF:147:ARG:NH2	2.35	0.59
51:AE:358:THR:HG22	51:AE:360:GLN:H	1.67	0.59
53:AG:71:THR:HG21	54:AI:252:GLN:CB	2.33	0.59
55:AJ:158:GLU:HG2	80:Ao:70:ILE:HG21	1.85	0.59
73:Ag:300:TRP:HH2	73:Ag:305:ILE:HD11	1.68	0.59
77:Ak:222:ASP:HA	77:Ak:225:VAL:HG22	1.84	0.59
4:BA:388:A:H2'	4:BA:389:A:H8	1.67	0.59
12:BL:178:ILE:O	12:BL:182:ASN:ND2	2.36	0.59
26:Bb:255:LEU:HD12	26:Bb:256:PRO:HD2	1.84	0.59
53:AG:122:GLN:NE2	53:AG:138:GLU:O	2.36	0.59
78:Am:47:THR:O	78:Am:50:SER:OG	2.20	0.59
27:Bc:159:TYR:CD2	27:Bc:186:LEU:HD11	2.38	0.59
45:Bw:242:PRO:HD3	45:Bw:423:PRO:HA	1.84	0.59
52:AF:15:ARG:NH2	70:Ad:178:GLU:OE1	2.36	0.59
57:AL:115:THR:HG21	57:AL:118:LEU:HD12	1.85	0.59
58:AN:113:HIS:O	58:AN:117:HIS:ND1	2.30	0.59
73:Ag:240:GLY:HA3	73:Ag:291:ASN:HD22	1.68	0.59
50:AC:89:ASP:O	50:AC:93:ARG:HG2	2.03	0.59
71:Ae:97:MET:HE1	71:Ae:131:TYR:HB3	1.85	0.59
73:Ag:154:ILE:HB	73:Ag:261:VAL:HG23	1.85	0.59
4:BA:157:A:OP2	87:B5:95:ARG:NH1	2.35	0.58
4:BA:788:A:O2'	17:BR:17:ARG:NH2	2.35	0.58
4:BA:1543:U:H2'	4:BA:1544:A:H8	1.68	0.58
29:Be:21:TRP:O	29:Be:32:LYS:NZ	2.36	0.58
38:Bn:72:GLU:N	38:Bn:72:GLU:OE2	2.33	0.58
45:Bw:63:VAL:HA	45:Bw:66:TRP:CD1	2.37	0.58
47:AA:330:A:H1'	56:AK:184:PRO:HB3	1.86	0.58
69:Ac:101:HIS:O	69:Ac:105:ILE:HG12	2.03	0.58
76:Aj:143:PRO:HB2	76:Aj:145:HIS:CE1	2.38	0.58
8:BF:94:ASP:OD2	15:BP:30:ASN:ND2	2.36	0.58
25:Ba:174:GLU:HA	25:Ba:296:SER:HB2	1.84	0.58
48:EL:74:LEU:HD11	48:FL:67:LEU:HD11	1.86	0.58
57:AL:63:LEU:HD11	57:AL:84:ARG:HB2	1.85	0.58
4:BA:1218:U:H5'	18:BS:173:ARG:HE	1.68	0.58
35:Bk:101:ALA:HB2	86:B4:44:LEU:HD12	1.85	0.58
67:Aa:338:GLU:HA	67:Aa:341:LYS:HG2	1.85	0.58
73:Ag:365:LEU:HD23	73:Ag:370:ALA:HB1	1.84	0.58
77:Ak:263:ASN:O	77:Ak:267:THR:HG23	2.04	0.58
4:BA:889:C:H41	4:BA:1082:U:H4'	1.68	0.58
4:BA:1383:G:OP1	12:BL:129:ARG:NH2	2.36	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
12:BL:113:GLU:O	12:BL:117:GLN:HG2	2.02	0.58
33:Bi:154:ASN:O	33:Bi:154:ASN:ND2	2.30	0.58
49:AB:133:HIS:HB3	68:Ab:34:ILE:HG12	1.85	0.58
54:AI:199:ARG:NH2	54:AI:205:GLU:OE1	2.35	0.58
59:AO:210:LEU:HD22	79:An:173:LEU:HG	1.84	0.58
4:BA:747:U:H3	29:Be:51:ILE:HG23	1.68	0.58
8:BF:123:GLY:HA3	8:BF:142:ARG:HG2	1.86	0.58
61:AQ:28:VAL:HG11	61:AQ:66:LEU:HD11	1.84	0.58
4:BA:108:A:N1	24:BY:40:ARG:NH2	2.51	0.58
4:BA:565:A:N6	4:BA:1314:U:OP1	2.35	0.58
4:BA:1326:A:O2'	4:BA:1328:G:OP2	2.20	0.58
40:Bp:80:HIS:O	46:Bx:36:ARG:NH1	2.37	0.58
47:AA:84:A:N6	47:AA:98:A:O2'	2.37	0.58
4:BA:197:U:H2'	4:BA:198:U:H6	1.69	0.58
4:BA:218:A:N6	4:BA:229:A:O4'	2.37	0.58
4:BA:777:A:N6	45:Bw:82:ILE:O	2.27	0.58
9:BI:102:VAL:HG22	9:BI:128:LEU:HD12	1.86	0.58
22:BW:101:LEU:HD22	22:BW:116:LYS:HG3	1.86	0.58
25:Ba:313:MET:HE1	25:Ba:353:GLN:HB2	1.85	0.58
47:AA:29:A:H2'	47:AA:30:A:C8	2.39	0.58
55:AJ:172:VAL:HB	77:Ak:157:ASP:HB3	1.85	0.58
9:BI:75:ARG:HH12	83:B1:100:ARG:NH2	2.01	0.58
40:Bp:89:ALA:O	40:Bp:93:HIS:ND1	2.24	0.58
67:Aa:204:ARG:NH2	67:Aa:210:GLU:OE2	2.37	0.58
76:Aj:82:ARG:NH2	76:Aj:132:GLU:O	2.37	0.58
4:BA:159:A:OP2	30:Bf:127:GLY:N	2.34	0.58
4:BA:234:G:N2	15:BP:33:SER:O	2.36	0.58
10:BJ:63:ARG:HH11	10:BJ:64:CYS:HB3	1.68	0.58
30:Bf:62:HIS:HE1	31:Bg:127:GLN:NE2	2.02	0.58
53:AG:37:TYR:HE2	54:AI:375:PRO:HD2	1.68	0.58
86:B4:52:LEU:HB2	86:B4:68:ARG:HG2	1.84	0.58
11:BK:25:ARG:HB3	11:BK:28:GLN:HB2	1.86	0.58
19:BT:57:PRO:HG2	71:Ae:130:TYR:HD2	1.69	0.58
47:AA:601:A:N6	75:Ai:63:GLY:HA3	2.19	0.58
71:Ae:236:LEU:HD22	71:Ae:240:GLU:HB3	1.86	0.58
83:B1:40:PRO:HG2	83:B1:45:PRO:HG3	1.85	0.58
30:Bf:105:GLN:HA	30:Bf:108:MET:HE2	1.86	0.57
33:Bi:170:PRO:HA	33:Bi:173:VAL:HG12	1.86	0.57
43:Bu:123:LYS:HA	43:Bu:126:ILE:HD12	1.86	0.57
47:AA:952:C:OP1	63:AU:52:ARG:NH2	2.37	0.57
64:AV:64:U:O4	64:AV:65:A:N6	2.36	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
73:Ag:173:ASN:O	73:Ag:176:ARG:NE	2.37	0.57
84:B2:93:ARG:NE	84:B2:150:GLU:OE1	2.37	0.57
4:BA:135:G:N1	4:BA:138:A:OP2	2.35	0.57
4:BA:843:C:O2'	6:BD:258:ILE:O	2.18	0.57
47:AA:3:A:O2'	47:AA:5:A:OP2	2.22	0.57
47:AA:272:A:H4'	47:AA:274:A:H4'	1.86	0.57
47:AA:799:A:H2'	47:AA:800:G:H8	1.69	0.57
61:AQ:12:TRP:HE3	61:AQ:65:LEU:HD11	1.69	0.57
68:Ab:71:ARG:HH12	72:Af:91:THR:HG22	1.68	0.57
71:Ae:142:TRP:HZ2	76:Aj:76:LEU:HD11	1.68	0.57
73:Ag:241:ILE:HD13	73:Ag:244:LYS:HE2	1.86	0.57
82:B0:114:VAL:O	82:B0:118:THR:HG23	2.04	0.57
4:BA:165:A:H2'	4:BA:166:A:C8	2.39	0.57
4:BA:1399:G:O2'	4:BA:1402:C:OP2	2.21	0.57
7:BE:121:LEU:HD22	7:BE:284:TYR:CE1	2.38	0.57
10:BJ:132:LYS:HE2	10:BJ:144:LEU:HD22	1.87	0.57
41:Bq:115:ARG:HH21	41:Bq:119:ARG:HH21	1.51	0.57
47:AA:441:U:H2'	47:AA:442:A:H8	1.70	0.57
60:AP:23:LEU:HD13	60:AP:32:TYR:HE1	1.68	0.57
70:Ad:169:THR:HG22	70:Ad:170:ARG:H	1.69	0.57
71:Ae:191:PHE:HA	71:Ae:196:ASN:HD22	1.70	0.57
4:BA:165:A:H2'	4:BA:166:A:H8	1.69	0.57
4:BA:677:C:OP1	4:BA:695:U:O2'	2.19	0.57
4:BA:1066:A:H2'	4:BA:1067:G:H8	1.69	0.57
16:BQ:98:HIS:HB3	16:BQ:151:VAL:HG12	1.85	0.57
45:Bw:168:SER:HG	45:Bw:232:HIS:HE2	1.51	0.57
47:AA:226:G:H2'	47:AA:227:A:H8	1.70	0.57
54:AI:198:SER:HB3	54:AI:246:ARG:HD2	1.87	0.57
71:Ae:122:ARG:HD3	71:Ae:159:TYR:HB2	1.86	0.57
88:B6:16:ILE:HD11	88:B6:61:LYS:HE2	1.85	0.57
8:BF:83:HIS:HD2	8:BF:85:ASP:HB2	1.70	0.57
22:BW:55:ILE:O	33:Bi:227:ARG:NH2	2.38	0.57
24:BY:110:PRO:HD3	33:Bi:166:GLU:HG2	1.85	0.57
50:AC:91:PHE:CZ	75:Ai:62:MET:HE2	2.39	0.57
56:AK:153:VAL:HG11	56:AK:164:ILE:HD11	1.87	0.57
77:Ak:216:LEU:O	77:Ak:220:ASN:ND2	2.34	0.57
80:Ao:338:PHE:HA	80:Ao:341:ILE:HG22	1.86	0.57
80:Ao:664:GLU:O	80:Ao:668:GLU:HG2	2.04	0.57
88:B6:42:THR:HG23	88:B6:57:VAL:HG12	1.87	0.57
4:BA:354:U:O2'	4:BA:1055:A:N1	2.35	0.57
22:BW:77:GLN:H	22:BW:171:HIS:CD2	2.22	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
43:Bu:52:SER:HB3	43:Bu:57:THR:HG21	1.87	0.57
45:Bw:240:GLN:NE2	45:Bw:348:PRO:O	2.37	0.57
62:AR:128:PRO:HA	62:AR:131:LEU:HB2	1.87	0.57
71:Ae:188:MET:O	71:Ae:192:ILE:HG12	2.05	0.57
73:Ag:126:TYR:CZ	73:Ag:343:LEU:HB2	2.40	0.57
81:Ap:105:CYS:HB2	81:Ap:108:CYS:CB	2.29	0.57
88:B6:16:ILE:HD12	88:B6:64:SER:HA	1.85	0.57
4:BA:222:G:N7	8:BF:170:ARG:NH2	2.46	0.57
26:Bb:231:GLU:OE2	26:Bb:300:THR:OG1	2.19	0.57
47:AA:115:A:OP1	59:AO:203:ARG:NH2	2.37	0.57
51:AE:395:PRO:HD2	51:AE:395:PRO:O	2.03	0.57
52:AF:37:ARG:HG2	52:AF:38:ASN:OD1	2.04	0.57
16:BQ:36:VAL:HG23	41:Bq:128:ARG:HH11	1.69	0.57
26:Bb:138:GLU:HG2	26:Bb:141:ARG:HH12	1.70	0.57
35:Bk:122:GLU:HB2	35:Bk:158:GLN:HG3	1.87	0.57
80:Ao:467:LEU:HB3	80:Ao:471:MET:HE1	1.85	0.57
4:BA:101:U:O3'	84:B2:230:ARG:NH2	2.38	0.57
4:BA:1218:U:H5'	18:BS:173:ARG:NE	2.20	0.57
4:BA:1379:C:O2	12:BL:98:LYS:NZ	2.38	0.57
8:BF:219:VAL:HG12	8:BF:265:THR:HG21	1.86	0.57
16:BQ:98:HIS:CD2	16:BQ:100:GLY:H	2.22	0.57
25:Ba:351:LEU:HD11	25:Ba:379:ASP:HA	1.87	0.57
25:Ba:361:THR:OG1	25:Ba:363:ASP:OD1	2.23	0.57
26:Bb:50:LYS:O	26:Bb:52:ARG:NH1	2.38	0.57
27:Bc:183:ASP:HB3	27:Bc:186:LEU:HD13	1.86	0.57
47:AA:471:U:OP2	68:Ab:50:ARG:NH2	2.38	0.57
55:AJ:123:SER:OG	55:AJ:124:VAL:N	2.38	0.57
70:Ad:168:ILE:HG12	70:Ad:176:ARG:HG3	1.87	0.57
87:B5:108:ASP:OD1	87:B5:109:VAL:N	2.38	0.57
4:BA:220:G:H1'	4:BA:231:C:H1'	1.86	0.57
4:BA:1230:U:H5'	4:BA:1231:U:H5'	1.87	0.57
4:BA:1269:G:N2	4:BA:1272:U:O2	2.31	0.57
8:BF:283:LEU:HB2	15:BP:125:ARG:HH12	1.68	0.57
11:BK:55:GLU:HA	11:BK:58:LYS:HE3	1.87	0.57
4:BA:402:A:OP1	35:Bk:48:TYR:N	2.37	0.56
10:BJ:123:MET:HB2	10:BJ:152:LEU:HD11	1.87	0.56
17:BR:108:LEU:HD11	17:BR:136:LEU:HD13	1.87	0.56
34:Bj:162:ARG:HE	34:Bj:171:TRP:HE3	1.52	0.56
73:Ag:207:TYR:OH	73:Ag:245:GLU:OE1	2.20	0.56
73:Ag:333:PHE:HA	73:Ag:336:LEU:HD23	1.87	0.56
4:BA:672:G:H21	4:BA:759:C:H5'	1.69	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:BA:925:G:N2	4:BA:964:A:OP2	2.35	0.56
4:BA:1380:A:H5''	12:BL:102:ARG:HH22	1.69	0.56
10:BJ:188:ARG:HH11	40:Bp:55:VAL:HB	1.70	0.56
28:Bd:129:HIS:CE1	86:B4:38:ARG:HH22	2.23	0.56
28:Bd:146:ALA:HB1	34:Bj:212:HIS:CD2	2.40	0.56
54:AI:292:GLY:HA3	54:AI:325:GLY:HA2	1.86	0.56
77:Ak:83:VAL:O	77:Ak:101:LYS:NZ	2.35	0.56
85:B3:78:ARG:O	85:B3:83:LYS:NZ	2.37	0.56
2:B8:94:VAL:N	4:BA:86:U:OP1	2.38	0.56
4:BA:843:C:N4	4:BA:871:A:O2'	2.38	0.56
4:BA:862:A:H5''	6:BD:213:THR:HG21	1.86	0.56
40:Bp:16:VAL:HG13	40:Bp:66:VAL:HG22	1.87	0.56
49:AB:168:ARG:HG2	54:AI:159:TYR:CE1	2.41	0.56
53:AG:61:GLU:HA	53:AG:64:TYR:CE1	2.40	0.56
54:AI:381:LYS:HA	54:AI:388:ALA:HB2	1.87	0.56
55:Aj:176:LYS:HB3	77:Ak:153:ILE:HB	1.87	0.56
63:AU:15:GLN:HG2	63:AU:16:GLU:HG2	1.87	0.56
80:Ao:547:LEU:HG	80:Ao:550:ALA:HB3	1.87	0.56
81:Ap:104:PRO:HB2	81:Ap:109:ARG:HB3	1.86	0.56
1:B7:82:ILE:HG22	1:B7:86:MET:HE2	1.88	0.56
4:BA:740:U:H2'	4:BA:741:A:H8	1.70	0.56
12:BL:128:ASN:ND2	12:BL:132:GLU:O	2.39	0.56
25:Ba:172:LYS:O	25:Ba:175:THR:OG1	2.23	0.56
47:AA:85:G:H2'	47:AA:86:A:H8	1.69	0.56
4:BA:920:C:H42	4:BA:924:G:H22	1.54	0.56
30:Bf:126:ARG:O	30:Bf:131:ARG:NH1	2.39	0.56
32:Bh:97:TYR:O	32:Bh:101:GLU:HG2	2.06	0.56
33:Bi:197:VAL:HG22	33:Bi:212:VAL:HG13	1.87	0.56
71:Ae:74:TYR:OH	71:Ae:409:GLU:O	2.22	0.56
4:BA:53:U:O4	4:BA:62:C:N3	2.38	0.56
10:BJ:110:LEU:HB3	11:BK:132:LEU:HD23	1.86	0.56
18:BS:39:VAL:HB	18:BS:42:GLU:HG2	1.87	0.56
23:BX:148:PHE:O	33:Bi:273:LYS:NZ	2.33	0.56
47:AA:590:C:H2'	47:AA:591:A:C8	2.41	0.56
54:AI:316:PHE:CD2	54:AI:370:LEU:HD21	2.41	0.56
71:Ae:312:ASP:HB2	71:Ae:314:GLN:HG3	1.86	0.56
4:BA:1281:A:O2'	4:BA:1283:U:O4	2.19	0.56
8:BF:179:THR:HG22	15:BP:13:LEU:HD11	1.86	0.56
23:BX:39:THR:HG22	23:BX:98:GLN:HG2	1.88	0.56
38:Bn:68:LYS:HE2	44:Bv:31:PRO:HB2	1.88	0.56
50:AC:108:LEU:HD23	50:AC:110:LEU:HD21	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
73:Ag:272:THR:HG23	73:Ag:315:LEU:HD13	1.88	0.56
83:B1:189:ASP:OD1	83:B1:190:ARG:N	2.39	0.56
4:BA:416:U:H2'	4:BA:417:G:C8	2.41	0.56
24:BY:48:PRO:HG3	84:B2:236:LEU:HD13	1.88	0.56
40:Bp:74:LEU:HD12	46:Bx:47:THR:HB	1.88	0.56
69:Ac:132:ARG:NH1	69:Ac:136:LEU:O	2.38	0.56
22:BW:67:ARG:HH12	33:Bi:228:PHE:HB3	1.70	0.56
26:Bb:220:ASN:HB3	26:Bb:232:TYR:HB2	1.87	0.56
28:Bd:87:ASP:OD1	35:Bk:153:HIS:HE1	1.87	0.56
32:Bh:59:ARG:NH1	32:Bh:181:SER:O	2.39	0.56
36:Bl:139:GLN:HB2	42:Bt:82:PHE:CE2	2.41	0.56
53:AG:48:LYS:NZ	54:AI:319:HIS:O	2.32	0.56
61:AQ:72:PRO:HB3	61:AQ:78:LYS:HG2	1.87	0.56
76:Aj:73:LEU:HD13	76:Aj:76:LEU:HD12	1.88	0.56
77:Ak:123:ARG:HG2	80:Ao:77:THR:HB	1.88	0.56
80:Ao:460:TYR:HA	80:Ao:463:LYS:HE2	1.88	0.56
4:BA:148:A:H4'	4:BA:149:A:H5''	1.88	0.56
8:BF:114:THR:O	8:BF:156:ARG:NH1	2.39	0.56
24:BY:65:LEU:HD11	24:BY:121:LYS:HE3	1.88	0.56
34:Bj:162:ARG:HB3	34:Bj:251:HIS:HB2	1.87	0.56
36:Bl:161:LEU:HD22	36:Bl:166:PHE:HD2	1.71	0.56
70:Ad:166:THR:O	70:Ad:176:ARG:NH2	2.38	0.56
80:Ao:603:ARG:HH21	80:Ao:636:LEU:HD12	1.70	0.56
4:BA:169:U:OP2	31:Bg:114:ARG:NH2	2.39	0.55
4:BA:175:C:H5''	13:BN:115:ASN:HB2	1.87	0.55
8:BF:276:HIS:O	36:Bl:90:ARG:NH2	2.35	0.55
33:Bi:118:SER:O	33:Bi:122:ILE:HG13	2.05	0.55
47:AA:58:U:P	76:Aj:53:ARG:HD2	2.46	0.55
47:AA:104:A:OP2	61:AQ:49:TYR:OH	2.17	0.55
55:AJ:145:LEU:HD13	55:AJ:148:LEU:HD11	1.89	0.55
84:B2:87:TRP:HB3	84:B2:92:LEU:HD21	1.89	0.55
5:BB:65:U:H2'	5:BB:66:A:C8	2.41	0.55
10:BJ:146:LEU:HD22	10:BJ:177:LEU:HB3	1.89	0.55
26:Bb:194:THR:OG1	26:Bb:197:GLU:OE1	2.23	0.55
47:AA:158:C:O3'	47:AA:172:A:N6	2.39	0.55
62:AR:141:TYR:HA	63:AU:32:MET:HE1	1.88	0.55
4:BA:225:C:OP1	15:BP:132:SER:OG	2.19	0.55
10:BJ:94:ARG:NH1	10:BJ:158:GLU:OE1	2.39	0.55
17:BR:153:LEU:HD23	27:Bc:315:ARG:HD2	1.88	0.55
26:Bb:233:VAL:O	26:Bb:296:ARG:NH1	2.40	0.55
27:Bc:75:VAL:HB	33:Bi:282:ILE:HB	1.88	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
28:Bd:143:GLN:HG2	35:Bk:165:THR:HG23	1.87	0.55
47:AA:527:G:O2'	47:AA:843:C:OP2	2.24	0.55
58:AN:71:ALA:HA	58:AN:74:GLU:CD	2.31	0.55
68:Ab:11:SER:HB2	68:Ab:44:PRO:HG2	1.89	0.55
77:Ak:189:LYS:O	77:Ak:193:ILE:HG12	2.07	0.55
80:Ao:80:ARG:NH1	80:Ao:487:SER:OG	2.38	0.55
15:BP:102:GLN:NE2	15:BP:106:ASP:OD2	2.39	0.55
23:BX:71:ARG:NH2	23:BX:73:GLN:OE1	2.39	0.55
47:AA:117:C:H42	59:AO:208:LYS:HE2	1.71	0.55
47:AA:882:A:H5'	71:Ae:108:SER:HB2	1.88	0.55
56:AK:99:ILE:HD11	56:AK:163:ALA:HB1	1.87	0.55
80:Ao:581:TYR:O	80:Ao:585:LEU:HG	2.06	0.55
4:BA:2:C:OP1	22:BW:48:LYS:N	2.34	0.55
4:BA:948:A:O2'	4:BA:1382:C:O2	2.24	0.55
4:BA:1506:A:H2'	4:BA:1507:G:C8	2.42	0.55
15:BP:140:LEU:HG	15:BP:156:VAL:HG11	1.89	0.55
21:BV:159:ARG:NH1	21:BV:161:GLU:OE2	2.36	0.55
49:AB:150:PHE:O	49:AB:154:ILE:HD12	2.07	0.55
33:Bi:211:GLN:HA	33:Bi:248:PHE:O	2.06	0.55
47:AA:57:A:O2'	47:AA:59:C:OP1	2.20	0.55
74:Ah:345:LEU:HD11	75:Ai:26:THR:HG21	1.87	0.55
80:Ao:488:VAL:HG12	80:Ao:489:PHE:CD2	2.42	0.55
4:BA:73:A:O2'	4:BA:224:A:O2'	2.21	0.55
19:BT:87:THR:HG21	19:BT:91:LYS:HD3	1.89	0.55
53:AG:112:THR:HG23	73:Ag:396:TYR:HD1	1.70	0.55
62:AR:125:TYR:HB3	63:AU:9:ALA:HB2	1.89	0.55
67:Aa:104:MET:HE3	67:Aa:290:PHE:CE2	2.41	0.55
67:Aa:184:SER:H	67:Aa:192:ARG:NH2	2.05	0.55
71:Ae:178:PHE:CG	76:Aj:72:PRO:HA	2.42	0.55
80:Ao:316:TRP:HA	80:Ao:319:ILE:HD12	1.89	0.55
85:B3:75:THR:HB	85:B3:83:LYS:HG2	1.89	0.55
1:B7:83:LEU:HD23	84:B2:210:PHE:CD2	2.42	0.55
25:Ba:228:ALA:HB3	25:Ba:292:TYR:HB2	1.89	0.55
35:Bk:195:LYS:HD2	35:Bk:198:PHE:HD2	1.71	0.55
51:AE:95:ALA:HB2	75:Ai:74:GLU:HB3	1.88	0.55
75:Ai:86:LYS:HG3	75:Ai:89:ARG:HH21	1.71	0.55
80:Ao:141:MET:HE3	80:Ao:141:MET:HA	1.88	0.55
4:BA:888:A:O2'	4:BA:891:A:N6	2.30	0.55
15:BP:234:LEU:HG	15:BP:238:ARG:HH21	1.72	0.55
21:BV:102:VAL:HB	21:BV:139:LEU:HB3	1.89	0.55
27:Bc:199:PHE:HB3	27:Bc:277:VAL:HG21	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
47:AA:313:A:O2'	47:AA:314:A:O4'	2.24	0.55
47:AA:525:U:H2'	47:AA:526:G:H8	1.71	0.55
73:Ag:318:PRO:HG2	73:Ag:321:ALA:HB2	1.88	0.55
81:Ap:213:LEU:HD23	81:Ap:217:ARG:HH21	1.72	0.55
4:BA:353:C:O2	4:BA:1267:A:O2'	2.24	0.55
4:BA:1569:A:N6	7:BE:152:VAL:O	2.40	0.55
10:BJ:225:GLN:HB3	10:BJ:226:PRO:HD3	1.89	0.55
29:Be:128:LEU:HD13	84:B2:156:ARG:HA	1.88	0.55
47:AA:374:C:OP2	78:Am:9:ARG:NH2	2.38	0.55
8:BF:83:HIS:CD2	8:BF:85:ASP:HB2	2.42	0.54
47:AA:617:C:H4'	51:AE:100:LYS:HG2	1.89	0.54
47:AA:692:U:H5''	75:Ai:89:ARG:HH12	1.72	0.54
70:Ad:74:PHE:HA	70:Ad:77:GLU:HG2	1.87	0.54
1:B7:88:LYS:NZ	4:BA:127:G:OP2	2.34	0.54
2:B8:163:THR:HG21	4:BA:76:G:H5'	1.89	0.54
4:BA:44:A:H5'	84:B2:200:ARG:HH12	1.72	0.54
22:BW:92:ILE:HA	22:BW:95:MET:HE3	1.87	0.54
26:Bb:161:LEU:HD13	26:Bb:271:LEU:HD11	1.89	0.54
27:Bc:273:PHE:HB2	27:Bc:300:VAL:HG12	1.89	0.54
50:AC:154:HIS:CG	77:Ak:77:PRO:HG3	2.43	0.54
52:AF:115:GLU:OE1	52:AF:116:LYS:HG3	2.07	0.54
18:BS:137:LEU:HB3	43:Bu:187:LEU:HD11	1.89	0.54
20:BU:73:THR:HG22	20:BU:83:TYR:HB2	1.89	0.54
47:AA:942:G:N2	47:AA:945:A:OP2	2.40	0.54
3:B9:82:ARG:NH1	4:BA:1512:U:OP2	2.40	0.54
4:BA:296:C:OP2	92:BA:3213:SPM:N1	2.41	0.54
4:BA:1518:A:O2'	46:Bx:99:MET:O	2.25	0.54
26:Bb:110:LEU:O	26:Bb:114:ARG:HG3	2.07	0.54
55:AJ:179:GLN:HG2	55:AJ:180:LEU:N	2.22	0.54
4:BA:52:A:O2'	4:BA:349:A:N3	2.38	0.54
4:BA:198:U:H2'	4:BA:199:A:C8	2.42	0.54
4:BA:571:A:N3	13:BN:111:MET:HG2	2.23	0.54
4:BA:644:G:OP1	87:B5:93:ARG:NH1	2.38	0.54
4:BA:733:A:OP1	6:BD:102:GLN:NE2	2.35	0.54
4:BA:935:C:H4'	4:BA:1426:G:H1'	1.89	0.54
13:BN:26:GLN:NE2	13:BN:147:GLN:OE1	2.40	0.54
16:BQ:89:ILE:HD11	16:BQ:176:LEU:HD22	1.89	0.54
31:Bg:27:GLN:OE1	31:Bg:28:ARG:NH2	2.40	0.54
34:Bj:56:PRO:HG3	34:Bj:237:LEU:HD11	1.89	0.54
34:Bj:245:GLN:HG3	34:Bj:248:LYS:HB3	1.88	0.54
40:Bp:33:PHE:HE1	40:Bp:64:VAL:HG11	1.73	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
71:Ae:192:ILE:HG13	71:Ae:224:TYR:CZ	2.42	0.54
76:Aj:167:PRO:HG2	76:Aj:170:LEU:HB3	1.88	0.54
4:BA:388:A:H2'	4:BA:389:A:C8	2.43	0.54
8:BF:221:LEU:HD22	8:BF:264:PRO:HB2	1.89	0.54
12:BL:193:HIS:NE2	47:AA:253:G:OP1	2.39	0.54
25:Ba:293:PHE:CD2	25:Ba:313:MET:HG2	2.42	0.54
39:Bo:78:VAL:HG21	85:B3:134:MET:HE1	1.89	0.54
51:AE:392:ARG:HG2	51:AE:394:ASP:H	1.71	0.54
62:AR:73:VAL:HG21	62:AR:107:GLU:HB3	1.89	0.54
3:B9:69:LEU:H	4:BA:560:A:H62	1.55	0.54
4:BA:885:G:N2	6:BD:281:GLY:O	2.34	0.54
10:BJ:49:ALA:HB1	42:Bt:32:LYS:HE2	1.89	0.54
10:BJ:169:LYS:HA	10:BJ:169:LYS:HE2	1.90	0.54
16:BQ:105:MET:HE3	16:BQ:122:TRP:HZ3	1.72	0.54
35:Bk:162:LEU:HD11	35:Bk:166:PHE:HD2	1.72	0.54
41:Bq:88:PRO:HG2	41:Bq:91:GLU:HG2	1.90	0.54
47:AA:799:A:H2'	47:AA:800:G:C8	2.43	0.54
51:AE:372:GLU:N	51:AE:382:ILE:O	2.36	0.54
71:Ae:167:TYR:O	71:Ae:171:ASN:ND2	2.40	0.54
73:Ag:312:THR:OG1	93:Ag:500:GTP:O2G	2.26	0.54
76:Aj:82:ARG:NH2	76:Aj:133:GLN:HB2	2.23	0.54
80:Ao:469:CYS:SG	80:Ao:498:ASP:HB3	2.48	0.54
16:BQ:228:LYS:HD3	85:B3:78:ARG:HH21	1.72	0.54
17:BR:55:TYR:CE1	19:BT:270:MET:HG3	2.43	0.54
17:BR:131:LEU:HD22	87:B5:134:THR:HG23	1.90	0.54
18:BS:96:HIS:HE2	18:BS:97:ARG:NH1	2.06	0.54
39:Bo:77:VAL:HG22	85:B3:99:VAL:HG21	1.89	0.54
47:AA:192:C:H2'	47:AA:193:A:H8	1.71	0.54
48:EL:74:LEU:HD11	48:FL:67:LEU:CD1	2.38	0.54
55:AJ:126:ILE:HG13	55:AJ:126:ILE:O	2.08	0.54
65:AX:2:A:H3'	65:AX:3:A:H8	1.73	0.54
73:Ag:122:ARG:NH1	73:Ag:336:LEU:O	2.41	0.54
74:Ah:273:GLU:O	74:Ah:276:LYS:HG3	2.08	0.54
81:Ap:215:ARG:HG3	81:Ap:218:ARG:HH21	1.72	0.54
4:BA:203:A:O2'	4:BA:239:C:O2	2.23	0.54
4:BA:702:A:OP1	84:B2:119:GLN:NE2	2.41	0.54
4:BA:713:A:H2'	4:BA:714:A:C8	2.43	0.54
4:BA:914:A:N1	47:AA:921:U:O2'	2.37	0.54
14:BO:37:ARG:HD2	14:BO:54:PRO:HG2	1.89	0.54
21:BV:143:ASP:OD2	32:Bh:310:ASN:ND2	2.30	0.54
28:Bd:192:TYR:OH	86:B4:60:GLN:NE2	2.40	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
45:Bw:51:LEU:HG	45:Bw:61:ARG:HE	1.72	0.54
47:AA:587:U:O2	47:AA:707:A:N6	2.41	0.54
55:AJ:120:LEU:HD21	58:AN:101:LYS:HE3	1.89	0.54
71:Ae:422:GLN:O	71:Ae:426:GLU:HG2	2.08	0.54
72:Af:115:ILE:HD13	72:Af:141:VAL:HG21	1.89	0.54
76:Aj:68:LEU:HD21	76:Aj:100:VAL:HG21	1.89	0.54
4:BA:150:A:N3	4:BA:198:U:O2'	2.37	0.54
26:Bb:138:GLU:O	26:Bb:142:THR:HG22	2.08	0.54
29:Be:44:ARG:HH12	45:Bw:157:ALA:HB3	1.73	0.54
45:Bw:289:GLY:HA3	45:Bw:329:TRP:HZ2	1.72	0.54
49:AB:137:ARG:HD3	68:Ab:34:ILE:HD11	1.90	0.54
52:AF:105:CYS:HB3	62:AR:69:CYS:HA	1.89	0.54
53:AG:204:LYS:O	53:AG:208:GLU:HG2	2.08	0.54
62:AR:87:PRO:HA	62:AR:126:LYS:HB3	1.90	0.54
73:Ag:47:ILE:HG23	73:Ag:65:GLN:HE22	1.73	0.54
84:B2:104:TRP:HE3	84:B2:144:LEU:HD22	1.73	0.54
8:BF:283:LEU:HB2	15:BP:125:ARG:NH1	2.23	0.53
10:BJ:85:GLU:O	10:BJ:89:VAL:HG23	2.08	0.53
10:BJ:105:SER:N	40:Bp:40:GLU:OE2	2.40	0.53
17:BR:130:CYS:SG	87:B5:156:THR:OG1	2.44	0.53
23:BX:27:GLN:HG2	84:B2:115:LEU:HB3	1.90	0.53
26:Bb:160:ASP:OD2	26:Bb:267:ARG:NH2	2.41	0.53
37:Bm:59:ARG:HG2	37:Bm:59:ARG:O	2.09	0.53
47:AA:759:U:H2'	47:AA:760:A:H8	1.73	0.53
50:AC:116:GLN:HB3	50:AC:150:PRO:HG2	1.90	0.53
1:B7:91:LYS:N	4:BA:118:U:OP1	2.37	0.53
4:BA:43:C:H1'	25:Ba:72:ARG:HH12	1.73	0.53
4:BA:439:A:OP1	16:BQ:73:ARG:NH1	2.37	0.53
14:BO:102:SER:OG	19:BT:158:GLN:NE2	2.36	0.53
28:Bd:68:LEU:HD21	35:Bk:211:LEU:HD12	1.89	0.53
34:Bj:58:LEU:HD12	34:Bj:153:LEU:HD22	1.89	0.53
59:AO:75:LEU:H	59:AO:107:ILE:HD11	1.73	0.53
80:Ao:420:GLN:HE21	80:Ao:459:PHE:HE2	1.56	0.53
80:Ao:461:TYR:CZ	80:Ao:488:VAL:HG11	2.43	0.53
4:BA:130:A:H2'	4:BA:131:A:O4'	2.08	0.53
27:Bc:61:LYS:HG2	33:Bi:275:PRO:HB3	1.89	0.53
32:Bh:80:GLN:O	32:Bh:210:ARG:NH2	2.42	0.53
45:Bw:397:GLU:HG2	45:Bw:398:THR:HG22	1.90	0.53
47:AA:61:G:N2	47:AA:198:C:O3'	2.42	0.53
47:AA:77:G:OP1	76:Aj:3:ARG:NH2	2.42	0.53
47:AA:198:C:H2'	47:AA:199:G:C8	2.44	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
51:AE:232:THR:HG21	51:AE:238:LYS:HD3	1.90	0.53
52:AF:37:ARG:HE	52:AF:69:TYR:HE1	1.55	0.53
67:Aa:197:ARG:CZ	67:Aa:203:LEU:HB2	2.39	0.53
71:Ae:222:SER:HA	71:Ae:225:VAL:HG22	1.90	0.53
71:Ae:265:LEU:HD11	71:Ae:303:MET:HE3	1.91	0.53
74:Ah:309:ASN:HA	77:Ak:165:ILE:HG22	1.91	0.53
77:Ak:191:LYS:NZ	77:Ak:237:LYS:O	2.41	0.53
80:Ao:396:ILE:HD13	80:Ao:425:VAL:HG21	1.90	0.53
4:BA:247:C:H2'	4:BA:248:A:H8	1.73	0.53
34:Bj:256:SER:OG	34:Bj:257:LYS:N	2.40	0.53
60:AP:64:LYS:HE3	67:Aa:178:TYR:CE1	2.44	0.53
70:Ad:131:GLN:O	70:Ad:135:GLN:HG3	2.09	0.53
71:Ae:242:ARG:NE	71:Ae:282:ALA:O	2.42	0.53
4:BA:661:U:O2'	17:BR:27:HIS:ND1	2.38	0.53
10:BJ:186:LEU:HD11	48:DL:77:LEU:HD11	1.91	0.53
11:BK:84:GLN:HG2	11:BK:124:LYS:HA	1.90	0.53
15:BP:14:ASP:OD1	15:BP:17:ARG:NH1	2.41	0.53
15:BP:44:ARG:HD3	15:BP:45:ARG:HG3	1.89	0.53
17:BR:32:LEU:HB3	17:BR:52:LEU:HD22	1.91	0.53
20:BU:114:LYS:HD3	30:Bf:45:CYS:SG	2.49	0.53
25:Ba:333:SER:HB3	25:Ba:363:ASP:HB3	1.89	0.53
31:Bg:133:PHE:HB3	32:Bh:259:ARG:HD2	1.90	0.53
50:AC:109:VAL:HB	50:AC:120:CYS:HB2	1.90	0.53
67:Aa:310:GLN:NE2	67:Aa:318:ASP:O	2.41	0.53
76:Aj:88:GLN:O	81:Ap:215:ARG:NH1	2.42	0.53
20:BU:71:ARG:NH1	22:BW:208:HIS:O	2.41	0.53
26:Bb:40:ILE:HD11	35:Bk:62:LEU:HD22	1.89	0.53
51:AE:242:ARG:HG3	51:AE:260:LYS:HB2	1.89	0.53
54:AI:111:LEU:HD22	77:Ak:114:LEU:HD23	1.90	0.53
55:AJ:145:LEU:HD11	55:AJ:156:TYR:CE1	2.44	0.53
77:Ak:188:ALA:O	77:Ak:235:SER:OG	2.27	0.53
83:B1:41:VAL:HG21	83:B1:83:GLU:HB3	1.91	0.53
4:BA:197:U:H2'	4:BA:198:U:C6	2.44	0.53
4:BA:719:C:O2'	25:Ba:112:ARG:HG3	2.08	0.53
8:BF:237:LEU:O	44:Bv:25:TYR:N	2.41	0.53
28:Bd:63:ALA:O	28:Bd:67:ARG:HG2	2.09	0.53
31:Bg:73:PRO:HB2	31:Bg:89:ILE:HG13	1.91	0.53
33:Bi:244:GLU:OE2	33:Bi:264:LYS:HE2	2.09	0.53
47:AA:287:C:N3	69:Ac:11:ARG:NH2	2.57	0.53
54:AI:321:LEU:HD13	54:AI:357:VAL:HG11	1.91	0.53
3:B9:76:CYS:SG	3:B9:89:CYS:HB2	2.48	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:BA:462:G:HO2'	4:BA:486:A:HO2'	1.57	0.53
4:BA:1418:G:N2	4:BA:1421:A:OP2	2.35	0.53
25:Ba:350:ARG:NH1	25:Ba:383:TYR:O	2.42	0.53
35:Bk:97:ALA:HB3	35:Bk:103:ALA:HB2	1.91	0.53
52:AF:3:ARG:NH1	52:AF:70:ALA:O	2.41	0.53
55:AJ:171:GLU:HG2	77:Ak:158:TYR:CE2	2.43	0.53
60:AP:20:ARG:NH2	60:AP:38:HIS:O	2.42	0.53
63:AU:72:ILE:O	63:AU:76:MET:HG2	2.09	0.53
82:B0:117:VAL:HA	82:B0:120:LEU:HD12	1.90	0.53
4:BA:96:U:N3	44:Bv:34:SER:O	2.42	0.53
10:BJ:218:THR:HA	10:BJ:221:MET:HE2	1.91	0.53
11:BK:64:ILE:HG21	11:BK:90:PHE:HE1	1.73	0.53
13:BN:3:SER:HA	32:Bh:309:GLU:HG3	1.92	0.53
33:Bi:122:ILE:HG22	33:Bi:126:ARG:NE	2.24	0.53
33:Bi:292:GLU:HG3	87:B5:179:ARG:HH22	1.74	0.53
34:Bj:44:PRO:HG2	34:Bj:227:LEU:HA	1.91	0.53
43:Bu:185:GLU:HG2	43:Bu:189:LYS:HE3	1.91	0.53
57:AL:106:HIS:ND1	57:AL:131:ASP:OD2	2.42	0.53
57:AL:137:LYS:NZ	57:AL:138:LYS:H	2.07	0.53
64:AV:9:C:H1'	64:AV:42:A:H2'	1.91	0.53
4:BA:49:A:H2'	4:BA:50:A:C8	2.44	0.52
4:BA:144:A:H2'	4:BA:145:A:C8	2.44	0.52
4:BA:698:U:H2'	4:BA:699:G:H8	1.74	0.52
17:BR:16:ARG:O	17:BR:25:ARG:NE	2.43	0.52
47:AA:17:G:H2'	47:AA:18:G:C8	2.44	0.52
50:AC:127:LEU:HB2	50:AC:132:PHE:CE1	2.44	0.52
4:BA:289:U:H2'	4:BA:290:A:C8	2.44	0.52
4:BA:1450:U:H2'	4:BA:1451:U:H6	1.75	0.52
15:BP:72:THR:HB	15:BP:77:ARG:HH21	1.74	0.52
47:AA:120:A:N6	47:AA:132:A:N1	2.57	0.52
64:AV:15:A:O2'	64:AV:16:A:H5''	2.10	0.52
77:Ak:71:VAL:HG12	77:Ak:72:TYR:HD1	1.74	0.52
4:BA:29:A:N3	4:BA:35:A:O2'	2.34	0.52
4:BA:139:G:C6	33:Bi:72:LEU:HD22	2.44	0.52
4:BA:1396:C:N4	12:BL:96:ASN:OD1	2.34	0.52
7:BE:157:LYS:HD2	7:BE:161:ILE:HG21	1.90	0.52
26:Bb:90:ILE:O	26:Bb:90:ILE:HG13	2.10	0.52
26:Bb:175:VAL:HG22	26:Bb:204:VAL:HG12	1.90	0.52
43:Bu:109:TRP:CD1	43:Bu:109:TRP:H	2.27	0.52
54:AI:235:GLY:N	54:AI:238:GLU:OE2	2.26	0.52
80:Ao:485:ILE:HD12	80:Ao:491:PRO:HG3	1.90	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:BA:918:U:OP1	64:AV:21:C:O2'	2.28	0.52
4:BA:1450:U:H2'	4:BA:1451:U:C6	2.45	0.52
13:BN:72:TRP:NE1	46:Bx:148:ASN:O	2.40	0.52
34:Bj:215:PHE:HB3	34:Bj:268:TYR:HE2	1.72	0.52
47:AA:50:U:H2'	47:AA:51:G:C8	2.39	0.52
47:AA:168:A:O2'	47:AA:169:A:O4'	2.22	0.52
47:AA:342:C:OP1	56:AK:96:ASN:ND2	2.42	0.52
49:AB:83:LEU:HB2	49:AB:117:LEU:HD13	1.92	0.52
50:AC:133:TYR:HA	50:AC:136:VAL:HG12	1.91	0.52
55:AJ:179:GLN:HA	77:Ak:149:PHE:HA	1.90	0.52
71:Ae:381:LYS:HE2	71:Ae:385:LEU:HD12	1.90	0.52
80:Ao:269:MET:HG2	80:Ao:277:GLN:HE21	1.74	0.52
80:Ao:496:MET:HB2	80:Ao:530:LEU:HD11	1.91	0.52
80:Ao:617:PHE:HD2	80:Ao:633:LEU:HD11	1.75	0.52
81:Ap:105:CYS:HB3	81:Ap:142:VAL:HA	1.91	0.52
81:Ap:224:LEU:O	81:Ap:225:ARG:NE	2.41	0.52
4:BA:699:G:H2'	4:BA:700:A:H8	1.74	0.52
4:BA:1194:U:H5''	26:Bb:364:ARG:HD2	1.91	0.52
4:BA:1425:A:H5''	12:BL:87:PRO:HG2	1.90	0.52
10:BJ:101:HIS:HB3	10:BJ:150:HIS:HB3	1.91	0.52
22:BW:121:GLU:O	22:BW:125:MET:HG3	2.09	0.52
37:Bm:76:LYS:HE2	37:Bm:83:VAL:HG21	1.91	0.52
47:AA:502:A:OP2	59:AO:202:ARG:NH2	2.42	0.52
49:AB:161:CYS:SG	49:AB:253:GLN:HA	2.50	0.52
64:AV:9:C:O2'	64:AV:10:A:N7	2.34	0.52
71:Ae:188:MET:HA	71:Ae:191:PHE:HD2	1.74	0.52
77:Ak:154:ASP:OD1	77:Ak:176:ARG:NH2	2.42	0.52
80:Ao:382:PHE:HB3	80:Ao:396:ILE:HD11	1.92	0.52
81:Ap:145:LYS:HA	81:Ap:148:LYS:HE2	1.91	0.52
83:B1:167:LEU:HD21	83:B1:202:GLU:HB3	1.90	0.52
4:BA:289:U:H2'	4:BA:290:A:H8	1.73	0.52
4:BA:881:C:H4'	4:BA:882:A:H5'	1.90	0.52
4:BA:1150:G:O2'	4:BA:1319:G:OP1	2.26	0.52
4:BA:1224:C:H2'	4:BA:1225:A:C8	2.45	0.52
12:BL:128:ASN:HD21	12:BL:132:GLU:HB2	1.75	0.52
17:BR:73:MET:HE1	17:BR:86:ILE:HG21	1.91	0.52
47:AA:29:A:O3'	57:AL:129:LYS:NZ	2.37	0.52
47:AA:546:U:OP1	53:AG:185:LYS:NZ	2.41	0.52
47:AA:939:A:H2'	47:AA:940:C:C6	2.45	0.52
54:AI:125:MET:HE1	77:Ak:85:LEU:HD12	1.92	0.52
60:AP:110:LEU:O	60:AP:114:ARG:HG2	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
76:Aj:24:GLN:HA	76:Aj:27:ARG:HE	1.74	0.52
4:BA:164:A:H4'	4:BA:165:A:C8	2.45	0.52
10:BJ:171:ILE:HD12	10:BJ:172:PRO:HD2	1.91	0.52
12:BL:91:ASN:O	12:BL:95:VAL:HB	2.10	0.52
25:Ba:297:ALA:HB3	25:Ba:303:ARG:HG2	1.92	0.52
27:Bc:285:THR:OG1	27:Bc:288:SER:O	2.20	0.52
35:Bk:48:TYR:HB3	35:Bk:51:LYS:HG2	1.90	0.52
47:AA:116:G:OP1	59:AO:207:LYS:NZ	2.38	0.52
51:AE:324:CYS:HB3	51:AE:329:ILE:HB	1.92	0.52
4:BA:1224:C:H2'	4:BA:1225:A:H8	1.74	0.52
20:BU:14:VAL:HB	20:BU:17:ARG:HH21	1.74	0.52
23:BX:144:VAL:HG22	33:Bi:232:MET:HE2	1.92	0.52
26:Bb:139:TRP:CD1	26:Bb:143:CYS:HG	2.28	0.52
47:AA:430:A:OP1	64:AV:35:C:O2'	2.22	0.52
47:AA:561:U:H3'	47:AA:562:A:C2	2.44	0.52
47:AA:620:C:H2'	47:AA:621:A:H8	1.74	0.52
47:AA:620:C:H4'	55:AJ:122:LYS:HG2	1.92	0.52
47:AA:829:G:H2'	47:AA:830:A:H8	1.75	0.52
55:AJ:157:LEU:O	55:AJ:161:GLN:HG3	2.09	0.52
60:AP:98:PRO:HB3	81:Ap:200:TYR:HE2	1.74	0.52
67:Aa:217:VAL:HG22	67:Aa:226:ILE:HD13	1.92	0.52
70:Ad:51:THR:HG21	81:Ap:228:SER:H	1.73	0.52
76:Aj:178:ARG:HB2	76:Aj:183:ASP:HB2	1.91	0.52
80:Ao:499:LEU:O	80:Ao:503:LEU:HG	2.09	0.52
86:B4:71:GLU:O	86:B4:73:ARG:N	2.38	0.52
4:BA:62:C:OP1	9:BI:76:ARG:NH1	2.43	0.52
7:BE:76:LYS:HG3	7:BE:170:LEU:HD21	1.92	0.52
15:BP:208:GLU:HB3	44:Bv:103:LEU:HD21	1.92	0.52
22:BW:80:TYR:HD2	22:BW:85:MET:HE2	1.74	0.52
31:Bg:28:ARG:HH12	32:Bh:71:GLU:CD	2.18	0.52
32:Bh:184:PHE:O	32:Bh:186:VAL:HG23	2.10	0.52
50:AC:38:ARG:HG2	50:AC:43:ARG:HH21	1.75	0.52
54:AI:74:ILE:O	54:AI:78:ILE:HG12	2.10	0.52
54:AI:263:TYR:HE1	54:AI:269:ALA:H	1.56	0.52
67:Aa:129:THR:HG22	67:Aa:130:GLN:H	1.75	0.52
72:Af:133:GLU:O	72:Af:136:GLN:NE2	2.43	0.52
73:Ag:264:VAL:HG21	73:Ag:307:LEU:HD13	1.91	0.52
83:B1:212:VAL:O	83:B1:216:ARG:HG2	2.09	0.52
86:B4:70:ARG:HD2	86:B4:71:GLU:HB2	1.91	0.52
4:BA:1076:A:O2'	83:B1:101:VAL:HG13	2.09	0.52
4:BA:1122:A:H2'	4:BA:1123:A:H8	1.74	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:BA:1380:A:H2'	4:BA:1381:A:C8	2.44	0.52
4:BA:1407:U:H5'	12:BL:87:PRO:HB3	1.90	0.52
6:BD:141:ALA:HB2	6:BD:154:ALA:HB2	1.92	0.52
7:BE:100:ILE:HD11	7:BE:177:LYS:HB2	1.92	0.52
15:BP:193:PHE:HE2	15:BP:200:ILE:HG12	1.75	0.52
32:Bh:148:LEU:HD12	32:Bh:149:PRO:HD2	1.90	0.52
37:Bm:66:ASP:OD1	37:Bm:69:SER:HB2	2.10	0.52
70:Ad:198:VAL:C	70:Ad:199:ARG:HD3	2.35	0.52
80:Ao:480:TRP:HE3	80:Ao:484:LEU:HD11	1.75	0.52
80:Ao:647:GLY:O	80:Ao:651:ARG:HG2	2.10	0.52
81:Ap:144:MET:O	81:Ap:148:LYS:HG2	2.10	0.52
86:B4:53:TYR:CG	86:B4:72:PRO:HB3	2.45	0.52
18:BS:39:VAL:HG12	18:BS:41:ASN:H	1.75	0.51
26:Bb:257:PRO:HB3	26:Bb:268:PHE:CZ	2.45	0.51
27:Bc:85:PRO:HG2	27:Bc:112:MET:HA	1.91	0.51
37:Bm:92:LEU:HD23	37:Bm:98:LYS:HG3	1.92	0.51
54:AI:88:LEU:HD11	54:AI:107:ALA:HB1	1.93	0.51
63:AU:25:THR:O	63:AU:29:ILE:HG12	2.10	0.51
4:BA:60:U:O2'	83:B1:96:LYS:O	2.28	0.51
7:BE:96:ARG:NH2	7:BE:301:ASP:OD1	2.43	0.51
8:BF:187:LEU:HD11	8:BF:261:LEU:HD12	1.91	0.51
22:BW:95:MET:SD	22:BW:99:GLN:HG2	2.50	0.51
31:Bg:35:ARG:HB2	42:Bt:101:TRP:HB2	1.91	0.51
47:AA:121:C:H2'	47:AA:122:C:C6	2.45	0.51
77:Ak:66:GLN:NE2	80:Ao:81:ASP:OD1	2.43	0.51
80:Ao:521:GLU:HG3	80:Ao:522:TYR:CD1	2.45	0.51
11:BK:35:LEU:HD21	11:BK:49:PHE:HE2	1.75	0.51
30:Bf:49:LEU:HD23	30:Bf:60:CYS:HB3	1.92	0.51
31:Bg:121:THR:HG21	31:Bg:124:PRO:HB3	1.93	0.51
47:AA:923:G:H1'	47:AA:944:A:H2	1.75	0.51
76:Aj:31:SER:OG	76:Aj:32:GLN:OE1	2.28	0.51
77:Ak:53:LYS:HZ2	77:Ak:55:LEU:HB2	1.76	0.51
4:BA:157:A:H2'	4:BA:158:A:H8	1.76	0.51
7:BE:198:PHE:HD1	7:BE:202:GLN:HE22	1.58	0.51
8:BF:231:VAL:HG13	44:Bv:26:ARG:HD2	1.92	0.51
27:Bc:41:GLN:HB3	27:Bc:258:ILE:HD12	1.92	0.51
27:Bc:278:SER:O	27:Bc:317:ARG:NH1	2.43	0.51
62:AR:66:CYS:HB2	62:AR:100:LEU:HA	1.92	0.51
4:BA:52:A:H4'	4:BA:245:A:H2	1.75	0.51
4:BA:412:C:O2'	82:B0:59:HIS:NE2	2.37	0.51
4:BA:789:A:O2'	7:BE:215:PHE:O	2.21	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:BA:1234:U:O2'	15:BP:76:LEU:O	2.24	0.51
4:BA:1425:A:O2'	64:AV:69:C:OP2	2.27	0.51
8:BF:183:ALA:HB2	15:BP:13:LEU:HD12	1.93	0.51
25:Ba:363:ASP:OD1	25:Ba:363:ASP:N	2.43	0.51
42:Bt:82:PHE:CD1	42:Bt:83:PRO:HD2	2.45	0.51
47:AA:56:U:H1'	47:AA:198:C:H1'	1.92	0.51
58:AN:62:ILE:HG21	74:Ah:341:HIS:HB3	1.93	0.51
71:Ae:213:PHE:HE1	71:Ae:221:LEU:HD12	1.76	0.51
76:Aj:98:THR:HG23	76:Aj:99:ARG:HG2	1.91	0.51
8:BF:49:ARG:NH1	8:BF:81:GLU:O	2.42	0.51
22:BW:67:ARG:NH2	33:Bi:228:PHE:O	2.34	0.51
26:Bb:329:ARG:HH11	26:Bb:329:ARG:HG3	1.76	0.51
34:Bj:147:THR:HG22	34:Bj:253:VAL:HG22	1.93	0.51
38:Bn:92:ILE:O	38:Bn:96:ARG:HG3	2.11	0.51
47:AA:11:G:O4'	47:AA:840:A:O2'	2.29	0.51
47:AA:164:C:H3'	47:AA:165:G:H8	1.75	0.51
47:AA:708:A:H3'	47:AA:709:A:H8	1.76	0.51
68:Ab:18:ASP:OD1	68:Ab:19:LEU:N	2.44	0.51
4:BA:1330:U:H5''	12:BL:94:LYS:HZ3	1.75	0.51
20:BU:70:ASN:O	20:BU:73:THR:OG1	2.25	0.51
24:BY:145:ARG:HH22	24:BY:157:PRO:HD3	1.75	0.51
45:Bw:132:ASP:OD1	45:Bw:134:THR:OG1	2.23	0.51
47:AA:235:A:N6	81:Ap:82:LYS:HG3	2.26	0.51
47:AA:590:C:H2'	47:AA:591:A:H8	1.75	0.51
53:AG:170:VAL:HG13	53:AG:237:ALA:HA	1.91	0.51
67:Aa:361:GLY:O	67:Aa:365:GLU:HG2	2.11	0.51
76:Aj:30:ASP:HA	76:Aj:33:ARG:HB3	1.92	0.51
80:Ao:579:LEU:HA	80:Ao:582:VAL:HG12	1.93	0.51
83:B1:180:ASP:HB3	83:B1:183:ARG:HD2	1.93	0.51
4:BA:53:U:H2'	4:BA:54:A:H8	1.76	0.51
4:BA:1017:U:HO2'	4:BA:1439:C:HO2'	1.58	0.51
21:BV:126:VAL:HG11	21:BV:132:ILE:HD11	1.92	0.51
38:Bn:53:MET:HE2	44:Bv:27:ALA:HA	1.92	0.51
47:AA:408:A:H4'	47:AA:949:G:N2	2.25	0.51
58:AN:61:THR:HG23	75:Ai:29:LYS:HD2	1.93	0.51
71:Ae:276:LEU:HB3	71:Ae:328:LYS:NZ	2.26	0.51
4:BA:849:U:OP2	6:BD:288:ARG:NH2	2.38	0.51
4:BA:1177:U:O2	16:BQ:63:ARG:NH2	2.44	0.51
7:BE:67:ASP:OD1	7:BE:68:GLU:N	2.44	0.51
16:BQ:172:VAL:HA	16:BQ:175:PHE:CE2	2.45	0.51
21:BV:133:ARG:HH12	21:BV:155:LYS:HG2	1.76	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
24:BY:172:ASP:OD2	29:Be:73:ASN:ND2	2.44	0.51
53:AG:146:HIS:O	53:AG:150:LYS:HG3	2.11	0.51
73:Ag:133:LYS:HG2	73:Ag:344:VAL:HG21	1.92	0.51
80:Ao:434:LEU:O	80:Ao:438:VAL:HG23	2.11	0.51
80:Ao:547:LEU:HD23	80:Ao:551:PHE:CE1	2.46	0.51
8:BF:94:ASP:O	8:BF:98:GLN:HG3	2.11	0.51
8:BF:253:MET:HE3	8:BF:259:LEU:HD22	1.93	0.51
43:Bu:105:ALA:HA	43:Bu:115:ARG:HD3	1.92	0.51
77:Ak:168:PRO:O	77:Ak:171:ARG:NH1	2.44	0.51
80:Ao:372:LEU:HD12	80:Ao:418:PHE:CD2	2.46	0.51
8:BF:71:GLY:H	8:BF:74:GLN:HE21	1.59	0.50
11:BK:78:PHE:CE2	11:BK:80:ILE:HD11	2.46	0.50
15:BP:191:VAL:HB	15:BP:192:PRO:HD3	1.93	0.50
56:AK:112:ALA:HB3	56:AK:137:ALA:HB2	1.92	0.50
60:AP:56:PRO:HA	60:AP:64:LYS:NZ	2.25	0.50
74:Ah:351:SER:OG	77:Ak:213:ARG:NH2	2.44	0.50
80:Ao:379:ILE:HD12	80:Ao:396:ILE:HG23	1.92	0.50
80:Ao:393:SER:HA	80:Ao:396:ILE:HD12	1.92	0.50
4:BA:607:A:O2'	20:BU:16:ASP:OD2	2.28	0.50
4:BA:1387:A:O2'	4:BA:1388:A:O5'	2.20	0.50
4:BA:1493:A:C6	17:BR:14:VAL:HG21	2.46	0.50
45:Bw:289:GLY:HA3	45:Bw:329:TRP:CZ2	2.46	0.50
53:AG:88:ASP:HB2	53:AG:91:ILE:HG12	1.92	0.50
54:AI:124:ILE:HD11	54:AI:126:LYS:HE3	1.93	0.50
80:Ao:500:LEU:HD11	80:Ao:534:ILE:HG22	1.93	0.50
80:Ao:666:LEU:O	80:Ao:670:THR:HG23	2.11	0.50
17:BR:149:LEU:HD23	27:Bc:312:LEU:HD13	1.93	0.50
27:Bc:100:SER:HA	27:Bc:126:THR:HG22	1.93	0.50
31:Bg:27:GLN:HB2	31:Bg:80:LEU:HD12	1.92	0.50
45:Bw:81:ARG:O	45:Bw:85:LYS:HB3	2.11	0.50
45:Bw:226:GLN:NE2	45:Bw:228:ASN:OD1	2.43	0.50
47:AA:226:G:H2'	47:AA:227:A:C8	2.46	0.50
47:AA:262:C:H2'	47:AA:263:G:H8	1.75	0.50
47:AA:552:A:H2'	47:AA:553:G:C8	2.46	0.50
47:AA:798:A:H2'	47:AA:799:A:H8	1.75	0.50
47:AA:948:U:H2'	47:AA:949:G:H8	1.75	0.50
53:AG:153:GLU:HA	53:AG:183:ALA:HB2	1.94	0.50
54:AI:78:ILE:HD13	54:AI:128:PRO:HG2	1.93	0.50
54:AI:162:MET:HE1	63:AU:86:GLY:HA2	1.93	0.50
55:AJ:78:VAL:HB	55:AJ:143:LEU:HB2	1.92	0.50
55:AJ:96:VAL:O	55:AJ:100:LYS:HG2	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
62:AR:77:ASN:O	62:AR:81:LEU:HG	2.12	0.50
74:Ah:344:GLU:HG3	74:Ah:345:LEU:HD12	1.93	0.50
80:Ao:419:PHE:HB2	80:Ao:451:ILE:HD11	1.94	0.50
4:BA:362:C:OP1	21:BV:186:ARG:NE	2.34	0.50
4:BA:1005:G:H5''	22:BW:110:LYS:HB2	1.92	0.50
15:BP:247:ILE:HG21	43:Bu:59:TRP:HD1	1.76	0.50
34:Bj:179:GLN:HG2	34:Bj:181:GLY:H	1.74	0.50
47:AA:182:A:H62	57:AL:55:ARG:HG2	1.76	0.50
47:AA:728:U:H2'	47:AA:729:A:C8	2.47	0.50
51:AE:376:GLU:HG2	67:Aa:115:PRO:HA	1.94	0.50
67:Aa:99:ASN:OD1	67:Aa:100:ILE:N	2.44	0.50
78:Am:68:LYS:H	78:Am:68:LYS:HD2	1.77	0.50
88:B6:20:MET:HG2	88:B6:58:GLU:HA	1.93	0.50
4:BA:465:A:N6	42:Bt:16:GLN:OE1	2.42	0.50
4:BA:699:G:H2'	4:BA:700:A:C8	2.47	0.50
4:BA:776:C:O2'	4:BA:778:A:N7	2.39	0.50
4:BA:937:C:H3'	4:BA:938:U:H5''	1.94	0.50
4:BA:1493:A:N6	17:BR:14:VAL:HG21	2.26	0.50
26:Bb:139:TRP:O	26:Bb:144:GLY:N	2.39	0.50
26:Bb:303:PHE:CE2	26:Bb:311:MET:HE1	2.47	0.50
28:Bd:182:ILE:H	86:B4:62:GLY:HA3	1.77	0.50
32:Bh:83:PHE:HE2	32:Bh:202:LEU:HB2	1.77	0.50
34:Bj:133:LEU:HD22	86:B4:64:THR:HG21	1.93	0.50
35:Bk:158:GLN:NE2	35:Bk:189:HIS:HD2	2.09	0.50
39:Bo:75:ARG:HE	85:B3:134:MET:HG3	1.76	0.50
55:AJ:55:PRO:HA	80:Ao:443:ASN:HD22	1.75	0.50
60:AP:86:PRO:HA	60:AP:89:LYS:HG3	1.92	0.50
71:Ae:70:LEU:HD13	71:Ae:182:TYR:OH	2.12	0.50
4:BA:863:U:O4	6:BD:247:ARG:NH2	2.45	0.50
4:BA:1196:G:OP2	26:Bb:368:ARG:NH1	2.45	0.50
4:BA:1387:A:H2'	4:BA:1388:A:C8	2.46	0.50
9:BI:108:ARG:NH1	9:BI:143:GLU:OE1	2.45	0.50
24:BY:138:THR:OG1	24:BY:142:GLU:O	2.26	0.50
25:Ba:207:ASN:HB3	45:Bw:151:LEU:HD11	1.94	0.50
33:Bi:121:SER:OG	33:Bi:199:CYS:SG	2.57	0.50
35:Bk:172:GLU:HA	35:Bk:175:HIS:NE2	2.26	0.50
46:Bx:72:ILE:CD1	46:Bx:115:ILE:HD11	2.40	0.50
47:AA:385:C:H2'	47:AA:386:U:H6	1.77	0.50
67:Aa:275:ILE:HG12	67:Aa:281:TYR:HD1	1.76	0.50
71:Ae:109:SER:HB3	71:Ae:145:ARG:HG2	1.93	0.50
71:Ae:153:ILE:HD11	71:Ae:187:LEU:HD21	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
73:Ag:275:ARG:HE	73:Ag:281:ILE:HG22	1.76	0.50
4:BA:247:C:H2'	4:BA:248:A:C8	2.47	0.50
8:BF:190:VAL:O	8:BF:262:THR:HA	2.12	0.50
12:BL:141:ARG:NH2	64:AV:70:C:OP2	2.44	0.50
47:AA:242:U:O4	47:AA:259:A:N7	2.44	0.50
53:AG:84:SER:HB3	53:AG:87:GLU:HG2	1.93	0.50
53:AG:158:LEU:HD12	53:AG:229:MET:HE3	1.93	0.50
73:Ag:290:HIS:HA	73:Ag:293:ARG:HD2	1.92	0.50
78:Am:64:GLU:OE2	78:Am:67:ARG:NH2	2.45	0.50
4:BA:158:A:N3	4:BA:1039:A:O2'	2.33	0.50
4:BA:428:A:N3	4:BA:600:C:O2'	2.43	0.50
4:BA:1034:G:OP1	13:BN:114:LYS:NZ	2.45	0.50
4:BA:1050:C:N4	4:BA:1323:U:H1'	2.27	0.50
8:BF:171:VAL:HG13	8:BF:273:LEU:HD22	1.93	0.50
22:BW:73:HIS:CD2	22:BW:118:VAL:HG13	2.42	0.50
22:BW:113:GLN:O	22:BW:117:GLU:HG3	2.11	0.50
30:Bf:68:PRO:HG2	30:Bf:71:HIS:CD2	2.44	0.50
47:AA:896:A:H2'	47:AA:897:C:C6	2.47	0.50
73:Ag:178:ASP:HB2	73:Ag:236:THR:HG21	1.94	0.50
4:BA:115:U:H2'	4:BA:131:A:H61	1.77	0.50
4:BA:220:G:N3	4:BA:231:C:O2'	2.45	0.50
4:BA:644:G:N2	4:BA:1008:A:OP2	2.36	0.50
10:BJ:63:ARG:HH21	46:Bx:102:ARG:HH21	1.59	0.50
15:BP:127:VAL:HG21	15:BP:138:VAL:HG21	1.94	0.50
20:BU:69:ILE:O	20:BU:73:THR:HG23	2.12	0.50
24:BY:167:PRO:HG3	29:Be:74:TYR:HE1	1.75	0.50
27:Bc:211:LEU:O	27:Bc:269:THR:OG1	2.28	0.50
45:Bw:63:VAL:HA	45:Bw:66:TRP:HD1	1.74	0.50
46:Bx:92:PHE:CZ	46:Bx:107:LEU:HD11	2.46	0.50
50:AC:93:ARG:NH1	50:AC:108:LEU:O	2.43	0.50
58:AN:41:ARG:NH1	75:Ai:48:THR:O	2.45	0.50
70:Ad:28:LYS:HD2	70:Ad:32:ASP:HB2	1.94	0.50
73:Ag:347:TYR:CZ	73:Ag:387:PRO:HB3	2.47	0.50
80:Ao:631:ILE:HG13	80:Ao:632:GLU:N	2.26	0.50
4:BA:16:A:H2'	4:BA:17:C:C6	2.47	0.49
4:BA:188:C:P	21:BV:181:ARG:HH21	2.34	0.49
4:BA:434:U:H2'	4:BA:435:C:C6	2.47	0.49
4:BA:468:A:C2	42:Bt:18:ILE:HD11	2.47	0.49
26:Bb:213:LEU:HB2	26:Bb:275:GLN:HG3	1.94	0.49
27:Bc:103:ALA:O	27:Bc:110:TRP:N	2.44	0.49
36:Bl:154:ASP:OD1	36:Bl:154:ASP:N	2.44	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
38:Bn:69:HIS:CD2	38:Bn:71:LYS:HG3	2.47	0.49
47:AA:386:U:H2'	47:AA:387:C:H6	1.77	0.49
54:AI:127:HIS:NE2	54:AI:129:GLU:OE1	2.45	0.49
77:Ak:239:GLU:OE1	77:Ak:241:TRP:NE1	2.45	0.49
77:Ak:255:TRP:HH2	77:Ak:264:ILE:HD12	1.76	0.49
80:Ao:372:LEU:HB2	80:Ao:415:ASP:CG	2.37	0.49
84:B2:157:LEU:HD21	84:B2:163:GLU:HA	1.94	0.49
4:BA:2:C:O2'	22:BW:147:ARG:O	2.17	0.49
4:BA:32:C:O2'	4:BA:36:A:N6	2.45	0.49
4:BA:645:U:H2'	4:BA:646:A:H8	1.77	0.49
5:BB:43:C:H5'	18:BS:121:ASN:ND2	2.27	0.49
7:BE:200:PRO:HG3	7:BE:332:LEU:HB3	1.93	0.49
12:BL:82:CYS:HB2	12:BL:98:LYS:HD3	1.94	0.49
15:BP:111:ASP:OD1	15:BP:111:ASP:N	2.45	0.49
16:BQ:46:GLU:HG3	16:BQ:47:LYS:HG2	1.94	0.49
25:Ba:140:ALA:HB2	25:Ba:416:ALA:HB2	1.93	0.49
33:Bi:71:ARG:HG3	33:Bi:71:ARG:O	2.12	0.49
4:BA:618:A:H62	38:Bn:41:ARG:NH1	2.10	0.49
13:BN:28:PRO:HG3	13:BN:67:PHE:CE1	2.47	0.49
25:Ba:287:CYS:SG	25:Ba:324:GLN:NE2	2.85	0.49
47:AA:386:U:O2'	52:AF:93:ILE:O	2.30	0.49
61:AQ:22:MET:HB3	61:AQ:25:THR:HB	1.94	0.49
76:Aj:166:TYR:O	81:Ap:199:TRP:NE1	2.39	0.49
80:Ao:259:ALA:HA	80:Ao:262:TYR:CE2	2.47	0.49
83:B1:86:VAL:HB	83:B1:105:TRP:HB2	1.94	0.49
4:BA:1238:A:H3'	4:BA:1239:A:H5''	1.94	0.49
8:BF:92:ARG:HB2	8:BF:176:VAL:HG21	1.95	0.49
8:BF:123:GLY:HA2	8:BF:141:ILE:HG12	1.94	0.49
12:BL:197:LYS:NZ	47:AA:9:U:OP1	2.42	0.49
13:BN:23:GLY:HA2	13:BN:31:LEU:HD11	1.93	0.49
14:BO:104:ASN:HD21	19:BT:158:GLN:NE2	2.06	0.49
14:BO:130:ARG:HD2	14:BO:134:PHE:CE1	2.46	0.49
16:BQ:55:ARG:HG3	16:BQ:56:VAL:HG23	1.94	0.49
19:BT:102:ARG:HH21	19:BT:106:LEU:HD22	1.77	0.49
20:BU:111:LYS:HB2	31:Bg:127:GLN:HE21	1.76	0.49
43:Bu:138:CYS:SG	43:Bu:139:SER:N	2.85	0.49
67:Aa:177:LYS:HB3	67:Aa:199:PRO:HD3	1.93	0.49
73:Ag:122:ARG:NH1	73:Ag:337:ASP:HB2	2.27	0.49
73:Ag:173:ASN:HB3	73:Ag:176:ARG:HE	1.78	0.49
76:Aj:52:MET:HA	76:Aj:55:TRP:NE1	2.27	0.49
82:B0:78:GLY:HA3	82:B0:132:HIS:CE1	2.47	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
82:B0:102:GLU:OE2	82:B0:132:HIS:NE2	2.39	0.49
87:B5:118:GLN:HB2	87:B5:121:ILE:HB	1.95	0.49
8:BF:249:ASN:ND2	38:Bn:55:GLY:O	2.45	0.49
15:BP:177:ALA:HA	15:BP:222:TYR:CD1	2.46	0.49
18:BS:110:TRP:NE1	18:BS:114:LYS:HD3	2.28	0.49
21:BV:146:THR:HB	30:Bf:60:CYS:SG	2.52	0.49
23:BX:133:GLN:O	23:BX:136:SER:OG	2.28	0.49
24:BY:126:THR:HB	24:BY:152:ARG:HH11	1.77	0.49
27:Bc:299:PRO:HD2	27:Bc:302:LEU:HB2	1.95	0.49
28:Bd:138:SER:HA	28:Bd:141:GLU:HG2	1.93	0.49
31:Bg:15:LEU:HD21	32:Bh:169:VAL:HA	1.95	0.49
34:Bj:173:LEU:HD22	34:Bj:233:PHE:CE2	2.47	0.49
46:Bx:124:ARG:HD3	46:Bx:156:PRO:HD3	1.94	0.49
49:AB:97:ARG:NE	49:AB:113:ASP:OD2	2.39	0.49
49:AB:232:THR:HA	68:Ab:44:PRO:HB3	1.94	0.49
51:AE:203:LEU:HD21	51:AE:268:PHE:HB3	1.93	0.49
53:AG:196:HIS:HB3	53:AG:199:VAL:HG23	1.94	0.49
53:AG:223:LYS:O	53:AG:227:HIS:CD2	2.66	0.49
54:AI:100:THR:O	54:AI:104:ILE:HG13	2.12	0.49
56:AK:113:SER:O	56:AK:117:GLU:HG2	2.12	0.49
56:AK:190:PRO:HG2	63:AU:45:TYR:HB2	1.94	0.49
67:Aa:364:ILE:O	67:Aa:368:LEU:HG	2.13	0.49
4:BA:410:U:O2'	4:BA:1169:A:N7	2.42	0.49
12:BL:72:ILE:HD11	12:BL:155:ARG:HG3	1.95	0.49
21:BV:115:GLU:HG2	31:Bg:19:LEU:HD11	1.95	0.49
28:Bd:128:GLU:OE2	86:B4:38:ARG:HD2	2.13	0.49
44:Bv:107:GLN:O	44:Bv:110:GLU:HG3	2.13	0.49
47:AA:262:C:H2'	47:AA:263:G:C8	2.47	0.49
47:AA:714:U:H2'	47:AA:715:G:H8	1.78	0.49
51:AE:141:TRP:CE2	75:Ai:70:LEU:HD11	2.47	0.49
77:Ak:76:ALA:O	77:Ak:112:ASN:ND2	2.39	0.49
77:Ak:250:MET:HG2	77:Ak:252:GLU:H	1.78	0.49
80:Ao:353:LYS:HB2	80:Ao:382:PHE:HE1	1.78	0.49
4:BA:230:G:H2'	4:BA:231:C:C6	2.47	0.49
4:BA:585:A:OP1	20:BU:58:ARG:NH1	2.45	0.49
5:BB:4:A:H3'	5:BB:5:A:H5''	1.95	0.49
10:BJ:127:PRO:HG2	10:BJ:130:VAL:HG22	1.95	0.49
11:BK:156:VAL:HG11	11:BK:159:LEU:HB2	1.94	0.49
17:BR:18:LEU:HB2	17:BR:25:ARG:HG3	1.94	0.49
24:BY:15:LEU:HD11	24:BY:26:PRO:HB3	1.94	0.49
34:Bj:186:GLN:OE1	34:Bj:186:GLN:N	2.45	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
47:AA:868:U:H3	47:AA:902:C:H42	1.59	0.49
53:AG:195:LYS:HB3	53:AG:204:LYS:HZ2	1.78	0.49
60:AP:27:THR:HG22	60:AP:28:ASN:HD22	1.77	0.49
67:Aa:224:ARG:NH1	67:Aa:256:GLN:O	2.45	0.49
71:Ae:149:ILE:HG21	71:Ae:179:PRO:HG3	1.94	0.49
71:Ae:370:GLN:HB3	71:Ae:374:ARG:NH1	2.24	0.49
73:Ag:80:HIS:HB3	73:Ag:155:PRO:HG3	1.93	0.49
76:Aj:61:GLU:HG3	76:Aj:83:LYS:CD	2.42	0.49
4:BA:199:A:H2'	4:BA:200:A:H8	1.78	0.49
4:BA:1131:U:H2'	4:BA:1132:G:H8	1.78	0.49
6:BD:83:GLY:HA2	6:BD:96:ILE:HG22	1.95	0.49
18:BS:54:ARG:HH22	18:BS:141:ILE:HG12	1.76	0.49
18:BS:125:CYS:SG	18:BS:157:SER:HB2	2.53	0.49
25:Ba:117:VAL:HG23	25:Ba:307:ASP:OD1	2.12	0.49
26:Bb:237:VAL:HG13	26:Bb:240:ILE:HD11	1.95	0.49
47:AA:621:A:H5''	55:Aj:130:HIS:CD2	2.48	0.49
54:AI:119:LYS:HA	54:AI:122:ARG:HH11	1.78	0.49
69:Ac:48:VAL:HA	69:Ac:52:ILE:HG12	1.93	0.49
71:Ae:80:TRP:HE3	71:Ae:83:ARG:HD2	1.78	0.49
72:Af:81:PHE:HA	72:Af:84:MET:HE3	1.95	0.49
73:Ag:162:LYS:HD3	73:Ag:315:LEU:HD21	1.94	0.49
1:B7:60:ASN:ND2	29:Be:25:ARG:O	2.46	0.49
1:B7:73:LEU:HD13	84:B2:174:ILE:HD11	1.95	0.49
10:BJ:89:VAL:HG22	10:BJ:124:LYS:NZ	2.27	0.49
14:BO:55:PRO:HB3	14:BO:77:ILE:HG12	1.93	0.49
18:BS:54:ARG:HD3	26:Bb:154:TYR:CE2	2.48	0.49
32:Bh:267:LEU:O	32:Bh:267:LEU:HD12	2.12	0.49
33:Bi:90:PRO:HB2	33:Bi:269:TRP:HH2	1.78	0.49
45:Bw:198:ARG:HD3	45:Bw:199:PRO:HD2	1.94	0.49
45:Bw:251:ILE:HG22	45:Bw:253:ILE:H	1.78	0.49
47:AA:85:G:H2'	47:AA:86:A:C8	2.47	0.49
47:AA:182:A:C6	57:AL:56:PRO:HD2	2.48	0.49
47:AA:583:A:C8	47:AA:584:C:H5	2.31	0.49
53:AG:61:GLU:HA	53:AG:64:TYR:HE1	1.77	0.49
73:Ag:66:HIS:HA	73:Ag:98:LEU:O	2.12	0.49
4:BA:187:G:H1'	4:BA:467:A:N6	2.28	0.49
4:BA:199:A:H5''	38:Bn:32:PHE:HD1	1.77	0.49
47:AA:945:A:OP1	79:An:141:HIS:NE2	2.36	0.49
56:AK:67:PRO:HG2	56:AK:82:PHE:CG	2.48	0.49
67:Aa:359:GLN:NE2	67:Aa:360:LYS:HE3	2.28	0.49
71:Ae:69:LEU:HB3	71:Ae:405:LEU:HD23	1.95	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
73:Ag:63:GLU:O	73:Ag:350:LYS:NZ	2.39	0.49
76:Aj:166:TYR:HD2	76:Aj:167:PRO:HD2	1.76	0.49
77:Ak:165:ILE:H	80:Ao:134:GLU:HG3	1.77	0.49
80:Ao:157:LEU:O	80:Ao:161:ILE:HG12	2.13	0.49
4:BA:108:A:OP2	32:Bh:31:ALA:N	2.46	0.48
8:BF:70:ARG:O	8:BF:202:TYR:OH	2.31	0.48
31:Bg:28:ARG:HD2	31:Bg:78:GLU:OE2	2.12	0.48
43:Bu:103:HIS:HE1	43:Bu:105:ALA:HB3	1.78	0.48
45:Bw:245:VAL:HG22	45:Bw:367:GLN:OE1	2.13	0.48
47:AA:122:C:H2'	47:AA:123:U:C6	2.48	0.48
47:AA:279:U:O2'	57:AL:54:GLY:O	2.31	0.48
47:AA:581:A:H61	47:AA:806:A:H3'	1.77	0.48
56:AK:165:LYS:NZ	56:AK:169:MET:SD	2.86	0.48
74:Ah:343:MET:HA	74:Ah:343:MET:HE3	1.95	0.48
80:Ao:373:ALA:HA	80:Ao:376:HIS:HB3	1.95	0.48
4:BA:648:C:H2'	4:BA:649:A:C8	2.48	0.48
4:BA:667:C:O2	45:Bw:271:GLN:NE2	2.46	0.48
7:BE:206:VAL:HG11	7:BE:289:VAL:HG13	1.93	0.48
9:BI:69:ARG:HH11	44:Bv:49:ARG:HH22	1.61	0.48
17:BR:117:ASP:OD2	17:BR:118:ARG:N	2.46	0.48
25:Ba:347:THR:HG22	25:Ba:349:GLY:H	1.78	0.48
32:Bh:67:ASP:O	32:Bh:71:GLU:HG2	2.13	0.48
33:Bi:122:ILE:HG22	33:Bi:126:ARG:CZ	2.42	0.48
47:AA:329:A:H1'	56:AK:185:HIS:CD2	2.48	0.48
50:AC:163:VAL:HG21	50:AC:167:ILE:HD12	1.94	0.48
56:AK:87:ILE:HD11	56:AK:150:ARG:HH21	1.79	0.48
67:Aa:195:VAL:HB	67:Aa:203:LEU:HD11	1.95	0.48
71:Ae:366:SER:HB3	71:Ae:367:LYS:HE2	1.95	0.48
80:Ao:536:MET:HB2	80:Ao:540:ARG:NH1	2.28	0.48
85:B3:129:PRO:HG3	85:B3:147:VAL:HG12	1.95	0.48
4:BA:644:G:O2'	4:BA:1007:G:O6	2.28	0.48
9:BI:120:ARG:HH22	83:B1:133:ASP:HA	1.78	0.48
25:Ba:198:LEU:HD13	25:Ba:418:TYR:HE2	1.79	0.48
37:Bm:68:GLN:O	37:Bm:72:GLU:HG2	2.13	0.48
46:Bx:89:LEU:HD21	46:Bx:118:CYS:HB3	1.95	0.48
48:CL:27:LYS:O	48:CL:30:GLN:HG3	2.13	0.48
51:AE:376:GLU:OE2	67:Aa:118:GLN:HG2	2.14	0.48
53:AG:187:MET:HG2	53:AG:219:VAL:HG21	1.94	0.48
70:Ad:55:VAL:O	70:Ad:58:GLU:HG3	2.14	0.48
70:Ad:64:ARG:HA	70:Ad:67:VAL:HG22	1.95	0.48
73:Ag:340:ILE:HG23	77:Ak:315:LYS:HD2	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
80:Ao:582:VAL:HA	80:Ao:585:LEU:HD12	1.95	0.48
4:BA:659:U:H5''	45:Bw:55:SER:HB2	1.95	0.48
7:BE:252:TRP:HB3	13:BN:84:PRO:HD3	1.95	0.48
16:BQ:36:VAL:HG23	41:Bq:128:ARG:HD2	1.95	0.48
23:BX:7:TYR:O	84:B2:185:GLN:NE2	2.46	0.48
34:Bj:55:ARG:HG3	34:Bj:238:LEU:HD23	1.96	0.48
34:Bj:177:ASP:O	34:Bj:190:ARG:NH2	2.44	0.48
35:Bk:201:ARG:O	35:Bk:205:GLU:HG2	2.13	0.48
45:Bw:363:PHE:O	45:Bw:390:THR:OG1	2.28	0.48
47:AA:524:U:H2'	47:AA:525:U:C6	2.49	0.48
48:CL:40:LEU:HB3	48:FL:72:ALA:HA	1.94	0.48
55:AJ:106:ILE:HD11	55:AJ:143:LEU:HB3	1.95	0.48
56:AK:96:ASN:OD1	56:AK:97:THR:N	2.46	0.48
67:Aa:104:MET:HE1	67:Aa:310:GLN:OE1	2.14	0.48
23:BX:144:VAL:HG21	33:Bi:276:ILE:HD12	1.96	0.48
27:Bc:153:ARG:NH2	27:Bc:256:ASP:O	2.36	0.48
27:Bc:282:ASN:HD21	27:Bc:289:LEU:HD12	1.76	0.48
27:Bc:313:LEU:O	27:Bc:317:ARG:HG2	2.12	0.48
28:Bd:67:ARG:HB3	28:Bd:71:ARG:NH1	2.28	0.48
30:Bf:62:HIS:CE1	31:Bg:127:GLN:NE2	2.81	0.48
31:Bg:35:ARG:HA	31:Bg:44:ARG:HD3	1.96	0.48
47:AA:235:A:OP1	51:AE:195:GLY:N	2.45	0.48
47:AA:878:A:H2'	47:AA:879:A:H8	1.78	0.48
49:AB:155:GLU:OE1	49:AB:167:THR:OG1	2.23	0.48
51:AE:198:TRP:HE3	51:AE:201:VAL:HG21	1.77	0.48
53:AG:176:ASP:OD1	53:AG:180:ARG:NH2	2.47	0.48
60:AP:23:LEU:HD13	60:AP:32:TYR:CE1	2.47	0.48
71:Ae:203:VAL:HA	71:Ae:206:GLU:OE2	2.13	0.48
76:Aj:98:THR:HG22	76:Aj:116:GLY:HA2	1.95	0.48
48:GL:75:THR:HG22	48:GL:78:GLU:HG2	1.96	0.48
4:BA:713:A:H2'	4:BA:714:A:H8	1.77	0.48
4:BA:971:A:O3'	79:An:163:ARG:NH2	2.46	0.48
32:Bh:251:SER:OG	32:Bh:252:ALA:N	2.46	0.48
47:AA:238:C:H2'	47:AA:239:U:C6	2.48	0.48
47:AA:551:G:N2	47:AA:778:G:O2'	2.46	0.48
71:Ae:434:GLU:O	71:Ae:438:ARG:HG2	2.13	0.48
73:Ag:373:GLU:HA	73:Ag:376:LYS:HE2	1.94	0.48
76:Aj:67:LEU:HD13	76:Aj:141:LEU:HB3	1.96	0.48
80:Ao:259:ALA:HA	80:Ao:262:TYR:CZ	2.49	0.48
80:Ao:644:VAL:O	80:Ao:644:VAL:HG12	2.13	0.48
4:BA:723:A:H3'	4:BA:724:A:H5''	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:BA:745:A:H3'	4:BA:746:U:H5''	1.94	0.48
4:BA:1323:U:H5'	4:BA:1324:C:OP2	2.14	0.48
16:BQ:102:PHE:HA	16:BQ:105:MET:HE2	1.96	0.48
27:Bc:102:LEU:HD21	27:Bc:109:PRO:HB3	1.96	0.48
27:Bc:127:PHE:HB3	27:Bc:302:LEU:HD11	1.95	0.48
45:Bw:144:LEU:HD12	45:Bw:170:PHE:CE1	2.49	0.48
47:AA:156:C:H2'	47:AA:157:A:C8	2.49	0.48
48:DL:75:THR:HG22	48:DL:78:GLU:HG2	1.96	0.48
49:AB:194:PHE:HZ	49:AB:207:ALA:HB3	1.77	0.48
54:AI:263:TYR:CD1	54:AI:268:MET:HA	2.48	0.48
59:AO:137:ILE:O	59:AO:141:GLU:HG3	2.13	0.48
67:Aa:243:GLN:O	67:Aa:245:GLN:HG3	2.13	0.48
76:Aj:89:HIS:ND1	76:Aj:121:LYS:HD3	2.29	0.48
77:Ak:204:THR:HG22	77:Ak:206:VAL:HG23	1.96	0.48
88:B6:16:ILE:HD13	88:B6:36:ARG:HH21	1.77	0.48
4:BA:44:A:H5'	84:B2:200:ARG:NH1	2.29	0.48
4:BA:838:A:H1'	4:BA:933:A:H61	1.78	0.48
4:BA:1236:C:H2'	4:BA:1237:A:H8	1.79	0.48
22:BW:59:PRO:HB3	33:Bi:225:TYR:CD1	2.49	0.48
26:Bb:362:PRO:HB2	26:Bb:364:ARG:HG2	1.95	0.48
51:AE:289:THR:HG21	51:AE:308:GLN:H	1.77	0.48
67:Aa:295:TRP:HA	67:Aa:298:VAL:HG22	1.96	0.48
75:Ai:13:ARG:HG2	75:Ai:17:ARG:HH21	1.77	0.48
4:BA:474:A:O2'	4:BA:596:A:O4'	2.31	0.48
4:BA:510:C:OP1	11:BK:99:LYS:NZ	2.47	0.48
4:BA:1009:A:H2'	4:BA:1010:A:C8	2.48	0.48
4:BA:1046:A:OP1	7:BE:239:ARG:NE	2.45	0.48
9:BI:99:THR:OG1	9:BI:127:GLY:O	2.32	0.48
10:BJ:233:LEU:HG	48:GL:79:ILE:HD11	1.96	0.48
25:Ba:259:ILE:HG23	25:Ba:263:MET:HE3	1.96	0.48
26:Bb:174:HIS:HB2	26:Bb:205:THR:OG1	2.14	0.48
32:Bh:105:ARG:HB2	32:Bh:112:LYS:HD2	1.96	0.48
47:AA:216:A:N7	47:AA:273:A:O2'	2.39	0.48
47:AA:242:U:H3	47:AA:259:A:H62	1.60	0.48
47:AA:552:A:H2'	47:AA:553:G:H8	1.78	0.48
48:FL:75:THR:HG22	48:FL:78:GLU:HG2	1.96	0.48
76:Aj:52:MET:HA	76:Aj:55:TRP:CD1	2.49	0.48
77:Ak:55:LEU:HD12	77:Ak:56:PRO:HD2	1.95	0.48
48:HL:75:THR:HG22	48:HL:78:GLU:HG2	1.96	0.48
87:B5:155:GLU:HB3	87:B5:172:LYS:HD3	1.95	0.48
4:BA:615:G:H5''	31:Bg:36:ASP:HB3	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:BA:668:A:H4'	45:Bw:89:MET:HE2	1.96	0.48
10:BJ:128:ASN:O	10:BJ:132:LYS:HG2	2.14	0.48
77:Ak:53:LYS:NZ	77:Ak:55:LEU:HB2	2.27	0.48
77:Ak:223:TYR:O	77:Ak:227:LEU:HD23	2.13	0.48
83:B1:46:HIS:CE1	83:B1:62:VAL:HG22	2.48	0.48
2:B8:108:LYS:NZ	4:BA:77:U:OP2	2.39	0.47
4:BA:1001:A:O2'	4:BA:1561:A:N3	2.45	0.47
4:BA:1251:C:H5'	4:BA:1253:G:H5'	1.96	0.47
7:BE:183:GLU:OE2	7:BE:183:GLU:N	2.46	0.47
8:BF:83:HIS:HD2	8:BF:85:ASP:H	1.61	0.47
8:BF:257:GLN:HG2	15:BP:21:ARG:NE	2.29	0.47
10:BJ:236:TYR:OH	48:HL:86:LEU:O	2.30	0.47
17:BR:150:LEU:HD23	17:BR:153:LEU:HD12	1.96	0.47
26:Bb:179:VAL:HB	26:Bb:183:ASP:HB3	1.96	0.47
34:Bj:45:TRP:HB3	34:Bj:230:LYS:HZ2	1.79	0.47
34:Bj:196:SER:OG	34:Bj:198:ASN:ND2	2.47	0.47
47:AA:104:A:OP1	61:AQ:10:ALA:N	2.42	0.47
47:AA:550:G:N1	47:AA:782:G:OP2	2.44	0.47
54:AI:366:ARG:HB3	54:AI:371:LEU:HD12	1.94	0.47
67:Aa:275:ILE:HG12	67:Aa:281:TYR:CD1	2.49	0.47
4:BA:301:U:H2'	4:BA:302:A:C8	2.50	0.47
4:BA:301:U:H2'	4:BA:302:A:H8	1.78	0.47
4:BA:311:A:H5'	6:BD:262:GLY:HA2	1.96	0.47
4:BA:639:U:H5'	22:BW:147:ARG:HH12	1.79	0.47
4:BA:1376:G:H1'	4:BA:1405:A:N3	2.28	0.47
5:BB:65:U:H2'	5:BB:66:A:H8	1.79	0.47
17:BR:106:ARG:HD3	17:BR:136:LEU:HD11	1.96	0.47
25:Ba:49:VAL:HG22	83:B1:160:ASP:OD2	2.14	0.47
47:AA:159:C:P	47:AA:172:A:H61	2.37	0.47
58:AN:90:ARG:HG2	58:AN:97:PRO:HA	1.96	0.47
76:Aj:31:SER:HA	76:Aj:111:HIS:CD2	2.48	0.47
4:BA:409:A:H61	35:Bk:48:TYR:HE2	1.61	0.47
4:BA:988:U:H4'	7:BE:230:THR:HG22	1.96	0.47
7:BE:52:HIS:NE2	27:Bc:301:HIS:O	2.47	0.47
7:BE:166:GLN:NE2	27:Bc:310:ASN:OD1	2.47	0.47
9:BI:55:ILE:HD12	9:BI:84:GLU:HB3	1.97	0.47
25:Ba:354:PHE:HB3	25:Ba:417:LEU:HD11	1.95	0.47
28:Bd:110:GLU:HA	28:Bd:113:ARG:HG2	1.95	0.47
31:Bg:144:ARG:HB2	31:Bg:147:GLU:HB2	1.95	0.47
46:Bx:79:HIS:HB3	46:Bx:184:ASN:HA	1.96	0.47
47:AA:365:A:H2'	47:AA:366:A:C8	2.49	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
47:AA:792:U:H4'	54:AI:389:ARG:HG3	1.96	0.47
47:AA:882:A:OP1	71:Ae:438:ARG:NH2	2.47	0.47
48:CL:25:PRO:HG2	48:DL:78:GLU:OE1	2.15	0.47
54:AI:273:GLY:HA3	54:AI:352:ALA:HB2	1.96	0.47
59:AO:115:LYS:HG3	59:AO:116:VAL:HG23	1.95	0.47
60:AP:38:HIS:CD2	60:AP:39:SER:H	2.31	0.47
68:Ab:115:GLU:O	68:Ab:119:VAL:HG22	2.14	0.47
2:B8:172:TYR:HB2	2:B8:175:ASP:HB2	1.95	0.47
2:B8:183:ARG:NH2	26:Bb:356:ARG:O	2.45	0.47
4:BA:199:A:H2'	4:BA:200:A:C8	2.49	0.47
4:BA:534:A:H5'	41:Bq:130:ASN:HB3	1.95	0.47
18:BS:90:ILE:HD13	18:BS:125:CYS:SG	2.54	0.47
24:BY:123:VAL:HG22	24:BY:130:PRO:HA	1.96	0.47
45:Bw:410:ASP:OD1	45:Bw:411:VAL:N	2.47	0.47
47:AA:368:A:C8	56:AK:185:HIS:ND1	2.83	0.47
47:AA:884:C:OP1	71:Ae:433:ARG:NH1	2.46	0.47
47:AA:903:C:H2'	47:AA:904:A:H8	1.79	0.47
53:AG:41:TYR:HB3	53:AG:75:LYS:HG3	1.96	0.47
54:AI:206:LEU:HD12	54:AI:219:TYR:HD1	1.80	0.47
60:AP:118:VAL:HG11	81:Ap:236:PRO:HB2	1.95	0.47
69:Ac:7:PHE:O	69:Ac:10:ARG:HG2	2.14	0.47
76:Aj:100:VAL:HG13	76:Aj:114:ALA:HB2	1.95	0.47
4:BA:554:U:O2'	46:Bx:102:ARG:NH1	2.36	0.47
4:BA:1060:C:H2'	4:BA:1061:U:H6	1.79	0.47
4:BA:1380:A:H2'	4:BA:1381:A:H8	1.80	0.47
6:BD:207:TYR:O	6:BD:209:ARG:HG3	2.13	0.47
8:BF:93:LEU:HD23	8:BF:96:LEU:HD12	1.95	0.47
10:BJ:54:ILE:HD12	16:BQ:211:ASN:HB3	1.96	0.47
26:Bb:185:MET:HE1	43:Bu:184:ARG:HB2	1.96	0.47
26:Bb:217:LEU:HD12	26:Bb:236:LEU:HD13	1.97	0.47
29:Be:85:THR:HG23	29:Be:88:GLN:H	1.79	0.47
32:Bh:93:VAL:HG12	32:Bh:180:LEU:HD12	1.97	0.47
34:Bj:51:LEU:HD12	34:Bj:188:ALA:HB1	1.96	0.47
46:Bx:117:GLU:HG3	46:Bx:180:LEU:HD12	1.96	0.47
47:AA:201:A:H2'	47:AA:202:A:C8	2.48	0.47
47:AA:568:G:N2	55:AJ:126:ILE:HG21	2.30	0.47
47:AA:883:C:H1'	47:AA:886:A:C6	2.49	0.47
51:AE:296:LEU:HD11	51:AE:303:ILE:HD12	1.97	0.47
59:AO:178:VAL:O	59:AO:182:ILE:HG12	2.15	0.47
60:AP:32:TYR:HB2	60:AP:54:TYR:HB3	1.95	0.47
61:AQ:50:PHE:HB3	61:AQ:77:VAL:HG22	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
62:AR:50:ASP:OD1	72:Af:86:ARG:NH1	2.46	0.47
62:AR:125:TYR:HE2	63:Au:4:HIS:HB3	1.78	0.47
62:AR:136:VAL:HG21	70:Ad:191:ILE:HG13	1.96	0.47
62:AR:141:TYR:O	63:Au:28:ARG:NH1	2.47	0.47
77:Ak:190:LYS:O	77:Ak:194:LYS:NZ	2.42	0.47
80:Ao:442:LEU:HD21	80:Ao:460:TYR:OH	2.15	0.47
4:BA:644:G:P	87:B5:93:ARG:HH12	2.36	0.47
4:BA:1000:A:H2'	4:BA:1001:A:C8	2.48	0.47
91:BA:3212:5GP:H8	91:BA:3212:5GP:OP3	2.14	0.47
10:BJ:229:LEU:HD11	48:HL:79:ILE:HD11	1.97	0.47
34:Bj:198:ASN:OD1	34:Bj:200:MET:HG2	2.15	0.47
34:Bj:235:LYS:HD2	34:Bj:275:PHE:HD1	1.79	0.47
47:AA:499:C:P	79:An:165:LYS:HE3	2.55	0.47
47:AA:814:A:OP2	54:Al:378:ARG:NH1	2.47	0.47
71:Ae:178:PHE:N	71:Ae:179:PRO:HD3	2.30	0.47
83:B1:26:ARG:HA	83:B1:26:ARG:HD2	1.69	0.47
4:BA:1401:U:H4'	7:BE:239:ARG:HH21	1.80	0.47
6:BD:198:GLU:HB3	6:BD:241:CYS:HB3	1.97	0.47
10:BJ:48:MET:O	16:BQ:250:ARG:NH1	2.42	0.47
10:BJ:198:PRO:HB3	10:BJ:202:LEU:HD23	1.95	0.47
11:BK:101:ALA:HB2	11:BK:109:ALA:HB2	1.97	0.47
16:BQ:105:MET:SD	16:BQ:183:LEU:HD21	2.55	0.47
18:BS:139:ALA:HA	26:Bb:189:HIS:O	2.15	0.47
25:Ba:99:TYR:HA	25:Ba:270:VAL:HG12	1.95	0.47
26:Bb:51:TYR:OH	82:B0:124:ALA:O	2.32	0.47
27:Bc:233:LYS:HD3	27:Bc:233:LYS:HA	1.64	0.47
32:Bh:245:LEU:HB3	32:Bh:250:ILE:HB	1.97	0.47
35:Bk:207:LEU:C	35:Bk:207:LEU:HD23	2.40	0.47
43:Bu:103:HIS:CE1	43:Bu:105:ALA:HB3	2.49	0.47
47:AA:25:U:H2'	47:AA:26:A:H8	1.79	0.47
49:AB:150:PHE:HZ	49:AB:242:PRO:HG3	1.79	0.47
52:AF:3:ARG:NH1	59:AO:98:MET:O	2.47	0.47
67:Aa:258:GLU:O	67:Aa:261:SER:OG	2.26	0.47
73:Ag:100:VAL:HG12	73:Ag:139:HIS:NE2	2.29	0.47
73:Ag:107:LEU:HD12	73:Ag:342:ILE:HD12	1.97	0.47
74:Ah:324:ILE:HG23	77:Ak:213:ARG:NH2	2.30	0.47
76:Aj:67:LEU:HB2	76:Aj:141:LEU:HD23	1.96	0.47
76:Aj:90:ASP:OD2	81:Ap:215:ARG:NH2	2.47	0.47
78:Am:102:LYS:O	78:Am:106:LYS:NZ	2.46	0.47
4:BA:1121:U:H2'	4:BA:1122:A:C8	2.50	0.47
4:BA:1414:C:N4	4:BA:1415:G:O6	2.48	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:BF:49:ARG:HG3	8:BF:263:LEU:HD22	1.97	0.47
18:BS:78:TRP:HB2	18:BS:174:GLU:OE1	2.15	0.47
23:BX:86:ARG:HH12	29:Be:18:MET:HB2	1.80	0.47
34:Bj:172:MET:HE1	34:Bj:175:GLN:H	1.80	0.47
47:AA:455:A:N7	47:AA:935:G:O2'	2.38	0.47
47:AA:815:U:H2'	47:AA:816:U:H6	1.79	0.47
47:AA:892:A:C5	76:Aj:21:LEU:HD21	2.50	0.47
67:Aa:307:LEU:O	67:Aa:311:ILE:HG12	2.15	0.47
71:Ae:246:ALA:HB2	71:Ae:289:LEU:HD11	1.96	0.47
77:Ak:68:TRP:N	77:Ak:69:PRO:HD2	2.30	0.47
4:BA:22:U:N3	84:B2:176:GLY:O	2.32	0.47
4:BA:469:C:OP2	85:B3:77:ARG:NH1	2.46	0.47
4:BA:482:G:OP1	39:Bo:23:ALA:N	2.47	0.47
4:BA:1072:A:H2'	4:BA:1073:A:C8	2.49	0.47
7:BE:202:GLN:HG2	7:BE:203:TYR:N	2.29	0.47
8:BF:220:ASP:OD1	8:BF:221:LEU:N	2.46	0.47
10:BJ:119:HIS:CE1	10:BJ:163:GLU:HB2	2.50	0.47
19:BT:275:THR:O	19:BT:279:GLU:HG3	2.15	0.47
23:BX:7:TYR:HE2	29:Be:56:PRO:HG3	1.79	0.47
41:Bq:112:PRO:HA	41:Bq:117:TYR:CD2	2.50	0.47
42:Bt:52:PRO:HA	85:B3:102:ASN:OD1	2.14	0.47
46:Bx:162:ILE:O	46:Bx:162:ILE:HG13	2.14	0.47
50:AC:165:LYS:HG3	50:AC:166:TYR:CD1	2.50	0.47
51:AE:408:TRP:NE1	51:AE:412:LYS:HD2	2.30	0.47
58:AN:34:MET:O	58:AN:38:VAL:HG23	2.15	0.47
58:AN:92:VAL:HB	58:AN:93:MET:HE2	1.97	0.47
71:Ae:302:VAL:HA	71:Ae:305:LYS:HE3	1.96	0.47
4:BA:792:G:H4'	7:BE:235:LYS:HB3	1.97	0.47
6:BD:144:ALA:HB2	6:BD:149:LYS:HG2	1.96	0.47
15:BP:207:PRO:O	15:BP:211:VAL:HG23	2.15	0.47
19:BT:152:ARG:HD3	19:BT:161:GLU:HG3	1.97	0.47
34:Bj:173:LEU:HG	34:Bj:174:PRO:HD2	1.97	0.47
47:AA:13:U:H2'	47:AA:14:C:C6	2.49	0.47
47:AA:468:A:OP2	49:AB:99:ARG:NH2	2.46	0.47
47:AA:578:G:H5''	47:AA:579:A:H2'	1.96	0.47
47:AA:843:C:O2	47:AA:928:A:N6	2.47	0.47
47:AA:939:A:H2'	47:AA:940:C:H6	1.80	0.47
53:AG:168:TYR:HB3	53:AG:236:LEU:HD23	1.97	0.47
54:AI:127:HIS:CD2	54:AI:129:GLU:HB2	2.50	0.47
55:AJ:104:ILE:HG21	55:AJ:145:LEU:HD22	1.96	0.47
67:Aa:344:ALA:HB3	67:Aa:349:LEU:HD23	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:BA:516:G:N3	4:BA:533:A:O2'	2.45	0.46
5:BB:44:U:H5''	28:Bd:119:LYS:HG2	1.97	0.46
11:BK:84:GLN:NE2	11:BK:85:PRO:HD2	2.30	0.46
33:Bi:121:SER:O	33:Bi:125:ILE:HG13	2.15	0.46
46:Bx:83:TYR:HB3	46:Bx:121:MET:HE3	1.97	0.46
47:AA:11:G:O6	47:AA:524:U:O4	2.33	0.46
47:AA:458:C:H5'	78:Am:22:LYS:HZ1	1.79	0.46
52:AF:60:ARG:HG2	52:AF:60:ARG:HH11	1.80	0.46
55:AJ:124:VAL:O	58:AN:108:ARG:NH2	2.45	0.46
64:AV:24:U:H2'	64:AV:25:C:C6	2.50	0.46
64:AV:50:U:O4	64:AV:56:C:N4	2.47	0.46
67:Aa:183:ILE:HA	67:Aa:192:ARG:HH21	1.80	0.46
67:Aa:213:ARG:HG2	67:Aa:214:MET:HE2	1.97	0.46
72:Af:150:LEU:HB2	72:Af:163:ILE:HG21	1.97	0.46
76:Aj:82:ARG:HH22	76:Aj:133:GLN:HB2	1.80	0.46
4:BA:290:A:O2'	4:BA:794:A:OP1	2.33	0.46
4:BA:471:U:OP1	85:B3:73:LYS:NZ	2.31	0.46
4:BA:744:A:OP1	4:BA:744:A:H4'	2.15	0.46
4:BA:806:U:H1'	7:BE:231:HIS:CE1	2.51	0.46
4:BA:1527:U:O2	4:BA:1527:U:H2'	2.16	0.46
11:BK:169:LYS:O	11:BK:173:LEU:HG	2.14	0.46
19:BT:79:GLU:OE2	19:BT:103:ARG:NH2	2.48	0.46
32:Bh:123:GLN:O	32:Bh:127:GLU:HG3	2.15	0.46
39:Bo:91:TRP:CD2	42:Bt:58:GLN:HG2	2.50	0.46
45:Bw:168:SER:OG	45:Bw:232:HIS:NE2	2.33	0.46
46:Bx:73:CYS:HB3	46:Bx:78:LYS:HE2	1.97	0.46
47:AA:10:U:N3	47:AA:13:U:OP2	2.33	0.46
47:AA:884:C:O4'	71:Ae:145:ARG:NH1	2.48	0.46
49:AB:141:ILE:O	49:AB:163:GLU:HB3	2.15	0.46
50:AC:98:GLY:HA3	75:Ai:71:TYR:CE1	2.50	0.46
67:Aa:233:LYS:O	67:Aa:237:LEU:HB2	2.15	0.46
71:Ae:113:SER:HA	71:Ae:116:VAL:HG12	1.95	0.46
73:Ag:271:THR:HG22	73:Ag:273:LEU:H	1.79	0.46
80:Ao:335:LEU:HD12	80:Ao:363:MET:HE3	1.95	0.46
83:B1:227:PHE:HD2	84:B2:158:LEU:HG	1.80	0.46
4:BA:392:U:H2'	4:BA:393:A:C8	2.45	0.46
4:BA:416:U:H1'	85:B3:105:SER:HB3	1.96	0.46
4:BA:484:A:OP1	42:Bt:37:ARG:NH1	2.49	0.46
4:BA:1020:C:H2'	4:BA:1021:C:C6	2.51	0.46
8:BF:171:VAL:HG12	8:BF:175:LYS:HE3	1.96	0.46
10:BJ:95:MET:HG3	10:BJ:181:ILE:CD1	2.46	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
13:BN:121:MET:HA	13:BN:124:ARG:HD2	1.97	0.46
15:BP:30:ASN:HB2	15:BP:33:SER:HB3	1.95	0.46
15:BP:117:ASP:OD2	15:BP:260:LYS:NZ	2.43	0.46
17:BR:65:LEU:HB3	17:BR:69:ASN:HD22	1.79	0.46
19:BT:195:PRO:O	19:BT:199:THR:HG23	2.15	0.46
45:Bw:111:LYS:HA	45:Bw:255:VAL:HG21	1.96	0.46
47:AA:568:G:H21	55:AJ:126:ILE:HG21	1.80	0.46
47:AA:826:C:H2'	47:AA:827:A:H8	1.81	0.46
49:AB:71:PHE:CE2	49:AB:262:LYS:HE2	2.50	0.46
53:AG:187:MET:HE1	53:AG:209:LEU:HD23	1.96	0.46
53:AG:223:LYS:O	53:AG:227:HIS:HD2	1.97	0.46
73:Ag:275:ARG:NH2	73:Ag:285:GLU:OE1	2.46	0.46
76:Aj:23:GLU:HB3	76:Aj:27:ARG:NH2	2.30	0.46
87:B5:125:TYR:O	87:B5:129:LYS:HG3	2.15	0.46
7:BE:105:GLY:HA3	19:BT:91:LYS:HD2	1.98	0.46
10:BJ:188:ARG:NH1	40:Bp:55:VAL:O	2.47	0.46
25:Ba:117:VAL:O	25:Ba:121:LEU:HG	2.16	0.46
30:Bf:115:MET:HG2	32:Bh:266:ALA:HB2	1.96	0.46
44:Bv:78:SER:OG	44:Bv:79:GLN:N	2.49	0.46
45:Bw:368:LEU:HD21	45:Bw:371:LEU:HD13	1.96	0.46
47:AA:96:U:H2'	47:AA:97:C:C6	2.50	0.46
47:AA:217:U:H2'	47:AA:218:A:H8	1.81	0.46
47:AA:670:A:H4'	53:AG:35:SER:HB2	1.97	0.46
47:AA:800:G:H2'	47:AA:801:A:H8	1.77	0.46
47:AA:903:C:H2'	47:AA:904:A:C8	2.51	0.46
49:AB:87:ARG:HB3	49:AB:90:LEU:HD12	1.97	0.46
51:AE:363:ALA:HB2	51:AE:370:VAL:HG23	1.96	0.46
58:AN:91:CYS:HB3	58:AN:96:ARG:H	1.80	0.46
70:Ad:76:SER:O	70:Ad:80:LYS:HD3	2.16	0.46
73:Ag:191:LYS:NZ	73:Ag:226:GLU:HG2	2.29	0.46
76:Aj:79:LEU:HD22	76:Aj:94:TYR:CD2	2.50	0.46
79:An:177:TRP:HB2	79:An:182:LEU:HB2	1.96	0.46
85:B3:68:ILE:HD11	85:B3:122:LEU:HD13	1.97	0.46
4:BA:233:A:H2'	4:BA:234:G:H8	1.81	0.46
4:BA:1287:G:H1	4:BA:1309:U:H3	1.64	0.46
4:BA:1416:G:H2'	4:BA:1417:A:H8	1.80	0.46
10:BJ:123:MET:C	10:BJ:124:LYS:HG3	2.41	0.46
14:BO:85:LEU:HD21	14:BO:134:PHE:CE2	2.51	0.46
14:BO:110:ASP:N	14:BO:110:ASP:OD1	2.46	0.46
17:BR:49:VAL:HG13	17:BR:124:ILE:HB	1.98	0.46
33:Bi:107:GLN:OE1	33:Bi:107:GLN:N	2.49	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:Bk:62:LEU:HD13	82:B0:124:ALA:HB1	1.98	0.46
49:AB:134:VAL:HG21	49:AB:191:LEU:HD12	1.98	0.46
49:AB:221:ILE:HA	49:AB:235:VAL:HG23	1.96	0.46
67:Aa:169:ILE:HG13	67:Aa:170:LEU:HG	1.98	0.46
72:Af:80:SER:O	72:Af:83:SER:OG	2.24	0.46
77:Ak:250:MET:HE2	77:Ak:252:GLU:HB2	1.96	0.46
80:Ao:302:GLU:OE2	80:Ao:306:LEU:HD23	2.15	0.46
80:Ao:375:TYR:O	80:Ao:379:ILE:HG12	2.15	0.46
80:Ao:423:MET:HB2	80:Ao:463:LYS:HE3	1.97	0.46
80:Ao:538:MET:HE1	80:Ao:585:LEU:HD11	1.97	0.46
81:Ap:223:ARG:HD2	81:Ap:223:ARG:N	2.31	0.46
4:BA:637:A:OP1	33:Bi:51:LYS:NZ	2.49	0.46
4:BA:740:U:H2'	4:BA:741:A:C8	2.50	0.46
4:BA:1156:A:O2'	82:B0:39:SER:N	2.49	0.46
4:BA:1518:A:H4'	46:Bx:94:ARG:NE	2.29	0.46
8:BF:53:LEU:HD13	8:BF:270:GLU:HG2	1.98	0.46
13:BN:67:PHE:HB3	13:BN:71:LYS:HB2	1.96	0.46
25:Ba:240:ALA:HB2	25:Ba:375:TRP:HH2	1.79	0.46
26:Bb:258:PHE:HA	26:Bb:327:VAL:HG13	1.97	0.46
40:Bp:40:GLU:OE2	40:Bp:43:ARG:NH1	2.48	0.46
47:AA:195:G:H5''	60:AP:20:ARG:HD2	1.96	0.46
47:AA:562:A:H8	47:AA:708:A:H61	1.64	0.46
47:AA:641:A:H2'	47:AA:642:G:C8	2.50	0.46
48:EL:67:LEU:O	48:EL:71:ILE:HG13	2.16	0.46
51:AE:217:ASP:OD1	51:AE:217:ASP:N	2.49	0.46
51:AE:243:VAL:HG11	51:AE:268:PHE:HE1	1.81	0.46
58:AN:65:LYS:HA	58:AN:68:GLN:HG3	1.97	0.46
71:Ae:408:CYS:O	71:Ae:411:GLU:HG2	2.15	0.46
73:Ag:237:ASP:O	73:Ag:241:ILE:HG12	2.16	0.46
75:Ai:86:LYS:HG2	75:Ai:91:LYS:O	2.16	0.46
76:Aj:129:ARG:HH12	76:Aj:204:PRO:HB2	1.80	0.46
48:GL:85:LEU:HA	48:GL:88:LYS:HG2	1.98	0.46
4:BA:31:A:N1	84:B2:200:ARG:N	2.58	0.46
4:BA:319:G:H2'	4:BA:320:U:C6	2.50	0.46
4:BA:344:U:H2'	4:BA:345:G:C8	2.51	0.46
18:BS:68:TRP:CD1	82:B0:78:GLY:HA2	2.51	0.46
22:BW:90:LYS:HD2	22:BW:93:ARG:HD2	1.96	0.46
24:BY:189:PRO:O	29:Be:93:THR:OG1	2.26	0.46
26:Bb:45:LEU:HD12	35:Bk:64:LYS:HE2	1.97	0.46
27:Bc:186:LEU:HG	27:Bc:189:TRP:HE1	1.81	0.46
33:Bi:124:LYS:HE2	33:Bi:208:ILE:HD13	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
45:Bw:186:ASN:HD22	45:Bw:420:LEU:HA	1.80	0.46
45:Bw:195:ASP:OD1	45:Bw:236:ARG:NH2	2.38	0.46
47:AA:798:A:H2'	47:AA:799:A:C8	2.50	0.46
48:DL:85:LEU:HA	48:DL:88:LYS:HG2	1.98	0.46
53:AG:71:THR:O	54:AI:254:LYS:NZ	2.48	0.46
73:Ag:275:ARG:HB2	73:Ag:277:ASP:OD1	2.15	0.46
76:Aj:52:MET:HA	76:Aj:55:TRP:CE2	2.51	0.46
77:Ak:78:PHE:CE1	77:Ak:107:LEU:HD13	2.50	0.46
80:Ao:541:ASP:OD1	80:Ao:542:GLN:N	2.48	0.46
81:Ap:105:CYS:CB	81:Ap:108:CYS:HB3	2.34	0.46
4:BA:1402:C:O2'	7:BE:233:GLN:HG3	2.16	0.46
16:BQ:101:HIS:O	16:BQ:105:MET:HG3	2.16	0.46
19:BT:275:THR:HA	19:BT:278:ILE:HD12	1.97	0.46
21:BV:150:LYS:HG3	30:Bf:58:ILE:HD11	1.96	0.46
24:BY:181:ASP:O	84:B2:95:LYS:NZ	2.40	0.46
27:Bc:91:HIS:ND1	33:Bi:281:MET:HB2	2.31	0.46
33:Bi:124:LYS:O	33:Bi:127:GLU:HG3	2.16	0.46
47:AA:855:C:H2'	47:AA:856:U:C6	2.50	0.46
48:EL:75:THR:HG22	48:EL:78:GLU:HG2	1.96	0.46
49:AB:81:ARG:NH2	49:AB:85:ASP:OD1	2.49	0.46
51:AE:243:VAL:HG11	51:AE:268:PHE:CE1	2.51	0.46
54:AI:127:HIS:HD2	54:AI:129:GLU:HB2	1.81	0.46
61:AQ:35:LEU:HB2	61:AQ:42:TYR:CE1	2.51	0.46
69:Ac:89:ASP:OD1	70:Ad:120:ARG:NH2	2.48	0.46
80:Ao:153:SER:O	80:Ao:157:LEU:HG	2.14	0.46
80:Ao:399:ILE:O	80:Ao:403:VAL:HG23	2.16	0.46
80:Ao:460:TYR:HD2	80:Ao:461:TYR:CD1	2.34	0.46
81:Ap:110:ASP:CG	81:Ap:113:LEU:HD23	2.40	0.46
48:HL:67:LEU:O	48:HL:71:ILE:HG13	2.16	0.46
4:BA:153:U:OP1	87:B5:87:ILE:HD13	2.15	0.46
4:BA:619:C:H2'	4:BA:620:A:H8	1.81	0.46
4:BA:709:A:H2'	4:BA:710:A:H8	1.80	0.46
4:BA:1527:U:H1'	46:Bx:154:TRP:HB3	1.98	0.46
13:BN:31:LEU:O	13:BN:34:MET:HG2	2.16	0.46
16:BQ:136:MET:HE3	16:BQ:136:MET:HB2	1.86	0.46
23:BX:138:ASP:HB3	23:BX:141:ARG:HG2	1.97	0.46
35:Bk:51:LYS:HD2	35:Bk:51:LYS:HA	1.72	0.46
35:Bk:127:MET:HB2	35:Bk:154:GLU:HB3	1.98	0.46
47:AA:114:A:N1	47:AA:136:C:O2'	2.44	0.46
47:AA:287:C:C2	69:Ac:2:PRO:HB3	2.51	0.46
47:AA:815:U:H2'	47:AA:816:U:C6	2.51	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
47:AA:956:G:H2'	47:AA:957:A:C8	2.51	0.46
51:AE:406:LEU:HD13	69:Ac:8:PRO:HB2	1.97	0.46
52:AF:76:GLU:O	52:AF:80:GLU:HG2	2.16	0.46
53:AG:189:ALA:O	53:AG:193:GLU:HG2	2.16	0.46
53:AG:199:VAL:HG11	53:AG:204:LYS:HD3	1.97	0.46
57:AL:32:THR:O	57:AL:36:MET:HG3	2.16	0.46
73:Ag:182:GLU:HA	73:Ag:185:THR:HG22	1.97	0.46
74:Ah:275:ALA:HA	74:Ah:278:LEU:HD12	1.98	0.46
74:Ah:311:ALA:O	77:Ak:166:ARG:NH2	2.47	0.46
77:Ak:68:TRP:HH2	77:Ak:84:PRO:HD2	1.80	0.46
77:Ak:286:LEU:HA	77:Ak:291:VAL:HG21	1.96	0.46
83:B1:147:LYS:HE3	83:B1:147:LYS:HB2	1.73	0.46
4:BA:33:A:C6	29:Be:30:PHE:HB2	2.51	0.46
19:BT:179:LEU:HD21	19:BT:215:VAL:HG22	1.96	0.46
27:Bc:212:PRO:HG3	27:Bc:268:ARG:NH1	2.31	0.46
47:AA:601:A:H2'	47:AA:602:A:C8	2.51	0.46
47:AA:729:A:O2'	73:Ag:316:PHE:N	2.49	0.46
47:AA:891:C:N4	47:AA:892:A:H62	2.15	0.46
47:AA:938:U:H2'	47:AA:939:A:C8	2.51	0.46
55:AJ:59:THR:HG22	55:AJ:61:PRO:HD3	1.97	0.46
48:FL:67:LEU:O	48:FL:71:ILE:HG13	2.16	0.46
73:Ag:54:ASP:N	73:Ag:54:ASP:OD1	2.49	0.46
73:Ag:144:CYS:SG	73:Ag:258:LEU:HD22	2.55	0.46
4:BA:48:U:O2'	4:BA:49:A:H4'	2.16	0.45
4:BA:253:A:N1	4:BA:334:C:H1'	2.30	0.45
4:BA:291:U:H4'	4:BA:793:A:H4'	1.97	0.45
4:BA:1390:G:H2'	4:BA:1391:U:C6	2.51	0.45
7:BE:244:SER:HB2	7:BE:251:VAL:HG12	1.99	0.45
23:BX:150:LEU:HD22	33:Bi:240:ARG:HE	1.81	0.45
25:Ba:124:THR:HG23	25:Ba:318:PHE:CG	2.51	0.45
35:Bk:127:MET:HE3	35:Bk:127:MET:HB3	1.82	0.45
47:AA:355:U:H2'	47:AA:356:A:H8	1.81	0.45
52:AF:54:HIS:CE1	52:AF:86:ILE:HA	2.51	0.45
53:AG:158:LEU:HD23	53:AG:171:PRO:HA	1.97	0.45
54:AI:158:TYR:CZ	54:AI:218:ASP:HB3	2.51	0.45
67:Aa:360:LYS:HB3	67:Aa:363:TYR:HD2	1.80	0.45
80:Ao:531:LYS:O	80:Ao:534:ILE:HG12	2.16	0.45
4:BA:170:U:O2	22:BW:205:THR:OG1	2.27	0.45
4:BA:1072:A:H2'	4:BA:1073:A:H8	1.81	0.45
4:BA:1338:G:H2'	4:BA:1339:A:H8	1.81	0.45
8:BF:91:PRO:HD3	15:BP:13:LEU:HD22	1.97	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
14:BO:42:ASP:HB3	14:BO:106:VAL:HG23	1.98	0.45
17:BR:38:ARG:NH2	17:BR:84:ASP:OD2	2.48	0.45
22:BW:189:THR:HG23	22:BW:192:ALA:H	1.81	0.45
33:Bi:169:PHE:CE2	33:Bi:173:VAL:HG11	2.52	0.45
34:Bj:214:LYS:HD3	34:Bj:230:LYS:HB3	1.97	0.45
40:Bp:19:GLN:NE2	40:Bp:65:ASP:OD2	2.50	0.45
44:Bv:48:PRO:HB2	44:Bv:50:TRP:CD1	2.51	0.45
47:AA:121:C:O2'	61:AQ:23:GLN:O	2.33	0.45
53:AG:204:LYS:HE2	53:AG:208:GLU:OE2	2.16	0.45
57:AL:137:LYS:HZ3	57:AL:138:LYS:H	1.63	0.45
64:AV:41:A:H2'	64:AV:42:A:O4'	2.16	0.45
67:Aa:245:GLN:O	67:Aa:249:VAL:HG23	2.16	0.45
68:Ab:19:LEU:HG	68:Ab:24:VAL:HB	1.98	0.45
71:Ae:427:ARG:O	71:Ae:431:ILE:HG12	2.16	0.45
73:Ag:221:LEU:HD11	73:Ag:242:VAL:HG22	1.99	0.45
75:Ai:18:LEU:HD12	77:Ak:195:LEU:HD11	1.98	0.45
83:B1:163:ARG:HD3	83:B1:205:GLY:H	1.80	0.45
4:BA:192:A:OP2	4:BA:1322:C:O2'	2.30	0.45
4:BA:216:C:OP2	36:Bl:111:ARG:NH2	2.48	0.45
4:BA:348:G:O6	38:Bn:113:ARG:NH1	2.43	0.45
4:BA:1557:C:O2'	4:BA:1558:U:O2	2.33	0.45
92:BA:3213:SPM:H62	92:BA:3213:SPM:H91	1.77	0.45
5:BB:31:C:H2'	5:BB:32:A:H8	1.80	0.45
7:BE:183:GLU:H	7:BE:183:GLU:CD	2.24	0.45
18:BS:122:VAL:HG22	18:BS:157:SER:HB3	1.99	0.45
22:BW:132:VAL:HG13	22:BW:178:GLU:OE1	2.16	0.45
23:BX:67:ALA:HB3	23:BX:98:GLN:HB2	1.97	0.45
28:Bd:61:ASP:OD1	28:Bd:61:ASP:N	2.47	0.45
29:Be:20:LYS:HB2	29:Be:20:LYS:HE2	1.80	0.45
32:Bh:121:ASP:OD1	32:Bh:122:ASN:N	2.49	0.45
34:Bj:173:LEU:O	34:Bj:175:GLN:HG3	2.17	0.45
49:AB:196:HIS:NE2	49:AB:239:ASP:O	2.49	0.45
53:AG:81:LYS:HD3	53:AG:81:LYS:HA	1.74	0.45
69:Ac:30:MET:HB3	69:Ac:64:ILE:HG12	1.97	0.45
71:Ae:426:GLU:O	71:Ae:430:LEU:HG	2.16	0.45
80:Ao:492:HIS:CG	80:Ao:493:SER:H	2.34	0.45
48:GL:67:LEU:O	48:GL:71:ILE:HG13	2.16	0.45
48:HL:85:LEU:HA	48:HL:88:LYS:HG2	1.98	0.45
4:BA:71:U:H3	4:BA:94:U:H3	1.64	0.45
4:BA:194:G:H2'	4:BA:195:G:C8	2.51	0.45
4:BA:1122:A:H2'	4:BA:1123:A:C8	2.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:BE:209:LYS:NZ	7:BE:265:ASP:OD1	2.39	0.45
11:BK:64:ILE:HG23	11:BK:89:TYR:CD1	2.52	0.45
21:BV:179:GLN:HG3	21:BV:184:TYR:HD2	1.81	0.45
27:Bc:66:HIS:HB2	27:Bc:76:PHE:HE1	1.81	0.45
27:Bc:76:PHE:CE2	33:Bi:281:MET:HG3	2.51	0.45
31:Bg:28:ARG:HG2	31:Bg:78:GLU:HB2	1.97	0.45
34:Bj:75:GLN:O	34:Bj:79:ILE:HG12	2.16	0.45
37:Bm:67:LEU:O	37:Bm:71:LEU:HD23	2.16	0.45
48:CL:39:THR:O	48:CL:43:ILE:HG12	2.16	0.45
49:AB:197:THR:OG1	49:AB:226:CYS:SG	2.63	0.45
50:AC:45:SER:HB2	50:AC:49:LYS:HD2	1.98	0.45
54:AI:223:ILE:O	54:AI:227:GLU:HG2	2.16	0.45
61:AQ:65:LEU:HD12	61:AQ:66:LEU:H	1.82	0.45
67:Aa:237:LEU:HG	67:Aa:241:TYR:CE2	2.51	0.45
71:Ae:280:LEU:O	71:Ae:283:VAL:HG22	2.17	0.45
72:Af:81:PHE:HE2	72:Af:85:LEU:HD22	1.80	0.45
74:Ah:367:TRP:HA	74:Ah:370:ASN:HD21	1.80	0.45
88:B6:47:ASP:O	88:B6:51:LYS:N	2.50	0.45
4:BA:860:G:H2'	4:BA:860:G:N3	2.31	0.45
4:BA:1003:C:C2	4:BA:1004:A:C8	3.05	0.45
4:BA:1357:C:H3'	4:BA:1358:G:H5''	1.98	0.45
6:BD:157:ASN:HB2	6:BD:179:GLU:OE1	2.17	0.45
18:BS:43:ALA:HB1	26:Bb:226:LEU:HD21	1.98	0.45
21:BV:165:ILE:HD11	21:BV:198:ARG:HB2	1.99	0.45
21:BV:168:THR:OG1	21:BV:169:GLU:N	2.49	0.45
31:Bg:35:ARG:NH2	42:Bt:102:SER:O	2.50	0.45
35:Bk:184:LEU:HD21	35:Bk:186:VAL:HG23	1.99	0.45
38:Bn:48:ASN:ND2	38:Bn:51:ARG:HH11	2.14	0.45
39:Bo:80:LEU:HD22	85:B3:101:LYS:HE3	1.98	0.45
46:Bx:70:CYS:SG	46:Bx:107:LEU:HA	2.57	0.45
47:AA:54:A:OP2	70:Ad:27:ARG:NH1	2.49	0.45
47:AA:204:U:H2'	47:AA:205:U:C6	2.52	0.45
47:AA:346:A:C6	65:AX:1:A:H1'	2.51	0.45
47:AA:432:A:H4'	47:AA:433:U:H5''	1.97	0.45
47:AA:789:C:H4'	47:AA:790:A:H5'	1.98	0.45
47:AA:878:A:H2'	47:AA:879:A:C8	2.52	0.45
49:AB:93:LYS:HE2	72:Af:149:GLU:HB3	1.99	0.45
67:Aa:344:ALA:HB1	67:Aa:348:HIS:HB3	1.98	0.45
71:Ae:392:SER:HB2	71:Ae:395:THR:HG22	1.97	0.45
76:Aj:197:ARG:HD2	76:Aj:197:ARG:HA	1.78	0.45
80:Ao:416:ASP:HB3	80:Ao:417:MET:HE2	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
81:Ap:60:ASP:OD1	81:Ap:121:LYS:NZ	2.44	0.45
4:BA:342:C:H2'	4:BA:343:C:H6	1.82	0.45
4:BA:838:A:H1'	4:BA:933:A:N6	2.30	0.45
4:BA:1239:A:H5'	4:BA:1240:A:H2	1.81	0.45
11:BK:84:GLN:HB2	11:BK:127:ASP:HB3	1.99	0.45
22:BW:75:ARG:HB3	22:BW:78:ILE:HD11	1.99	0.45
34:Bj:85:SER:HB3	34:Bj:88:GLU:HG2	1.98	0.45
36:Bl:72:TRP:NE1	36:Bl:92:HIS:O	2.37	0.45
44:Bv:61:PHE:CD1	44:Bv:65:GLY:HA2	2.51	0.45
47:AA:13:U:H2'	47:AA:14:C:H6	1.81	0.45
47:AA:88:C:H2'	47:AA:89:U:C6	2.52	0.45
47:AA:136:C:O2	59:AO:198:ARG:NH2	2.50	0.45
47:AA:168:A:O2'	47:AA:170:A:N1	2.49	0.45
48:DL:67:LEU:O	48:DL:71:ILE:HG13	2.16	0.45
50:AC:120:CYS:HB3	77:Ak:108:LEU:HD23	1.99	0.45
60:AP:102:MET:HE3	60:AP:106:ASN:ND2	2.22	0.45
67:Aa:184:SER:H	67:Aa:192:ARG:HH22	1.64	0.45
72:Af:144:ARG:HH21	72:Af:171:LEU:HD21	1.81	0.45
72:Af:144:ARG:NH2	72:Af:171:LEU:HD21	2.31	0.45
76:Aj:91:GLU:HB2	76:Aj:121:LYS:HD2	1.99	0.45
87:B5:132:LYS:O	87:B5:136:GLU:HG2	2.17	0.45
4:BA:1000:A:H2'	4:BA:1001:A:H8	1.81	0.45
7:BE:228:PRO:HG2	7:BE:233:GLN:NE2	2.32	0.45
11:BK:39:LEU:HD12	11:BK:44:VAL:HB	1.99	0.45
11:BK:118:TYR:CD1	41:Bq:96:LEU:HD21	2.52	0.45
21:BV:133:ARG:NH2	21:BV:155:LYS:O	2.50	0.45
24:BY:135:TRP:CE3	24:BY:143:ARG:NH1	2.85	0.45
25:Ba:118:LYS:NZ	25:Ba:254:GLU:OE1	2.37	0.45
27:Bc:43:ASP:O	27:Bc:47:LYS:HG2	2.17	0.45
47:AA:349:A:N3	47:AA:426:G:O2'	2.47	0.45
47:AA:956:G:H2'	47:AA:957:A:H8	1.80	0.45
55:AJ:84:ASP:OD1	55:AJ:84:ASP:N	2.45	0.45
56:AK:90:ILE:HD11	56:AK:163:ALA:CB	2.47	0.45
67:Aa:164:LEU:HD13	67:Aa:203:LEU:HG	1.99	0.45
71:Ae:121:SER:HA	71:Ae:159:TYR:CE2	2.45	0.45
73:Ag:82:LEU:HD12	73:Ag:83:PRO:HD2	1.98	0.45
85:B3:109:LYS:O	85:B3:113:VAL:HG22	2.17	0.45
4:BA:61:U:P	83:B1:100:ARG:HH21	2.38	0.45
4:BA:433:G:H22	4:BA:458:A:H2	1.65	0.45
4:BA:505:U:C2	4:BA:507:G:H5'	2.52	0.45
4:BA:1344:C:H2'	4:BA:1345:A:C8	2.52	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:BA:1378:U:H3	12:BL:82:CYS:HB3	1.82	0.45
6:BD:198:GLU:OE2	6:BD:202:GLY:N	2.48	0.45
7:BE:124:VAL:HG13	7:BE:127:CYS:SG	2.56	0.45
8:BF:99:VAL:O	8:BF:103:GLN:HG3	2.17	0.45
10:BJ:113:ARG:O	10:BJ:117:ARG:HG2	2.17	0.45
16:BQ:86:ARG:NH2	16:BQ:167:CYS:O	2.49	0.45
19:BT:188:LEU:HD12	19:BT:191:ARG:HH12	1.81	0.45
25:Ba:357:LEU:HD12	25:Ba:374:ALA:HB2	1.99	0.45
26:Bb:151:LEU:HD22	26:Bb:189:HIS:ND1	2.32	0.45
26:Bb:160:ASP:HA	26:Bb:163:HIS:CE1	2.52	0.45
28:Bd:121:TRP:CE3	34:Bj:70:MET:HG3	2.52	0.45
34:Bj:162:ARG:HB2	34:Bj:171:TRP:CZ3	2.52	0.45
43:Bu:75:ARG:NH1	43:Bu:108:ASP:OD1	2.50	0.45
45:Bw:87:GLN:HE22	45:Bw:273:GLU:HA	1.82	0.45
46:Bx:70:CYS:SG	46:Bx:108:CYS:N	2.88	0.45
47:AA:386:U:H2'	47:AA:387:C:C6	2.51	0.45
47:AA:584:C:O2	58:AN:86:ARG:NE	2.50	0.45
47:AA:661:U:O2'	54:AI:118:GLU:OE1	2.33	0.45
50:AC:99:THR:O	50:AC:101:PRO:HD3	2.17	0.45
52:AF:2:PRO:HB2	52:AF:96:HIS:CG	2.52	0.45
54:AI:313:GLN:O	54:AI:346:ARG:HG3	2.17	0.45
56:AK:65:TYR:CG	62:AR:131:LEU:HD21	2.51	0.45
73:Ag:123:TYR:CE1	73:Ag:340:ILE:HD12	2.52	0.45
76:Aj:110:ASP:O	76:Aj:111:HIS:ND1	2.50	0.45
4:BA:913:G:H21	4:BA:914:A:N6	2.15	0.45
4:BA:1264:C:OP2	4:BA:1409:C:O2'	2.35	0.45
12:BL:102:ARG:HD3	12:BL:132:GLU:OE2	2.16	0.45
20:BU:50:PHE:O	20:BU:54:THR:HG23	2.15	0.45
20:BU:111:LYS:HB2	31:Bg:127:GLN:NE2	2.32	0.45
25:Ba:209:LEU:O	25:Ba:223:HIS:HA	2.17	0.45
29:Be:121:GLN:NE2	29:Be:128:LEU:HD12	2.29	0.45
33:Bi:154:ASN:C	33:Bi:154:ASN:ND2	2.74	0.45
35:Bk:93:ILE:O	35:Bk:156:VAL:HA	2.17	0.45
35:Bk:111:HIS:CE1	35:Bk:121:VAL:HG11	2.52	0.45
40:Bp:63:CYS:SG	40:Bp:64:VAL:N	2.90	0.45
47:AA:154:C:H2'	47:AA:155:A:H8	1.81	0.45
47:AA:155:A:H2'	47:AA:156:C:C6	2.51	0.45
47:AA:525:U:H2'	47:AA:526:G:C8	2.50	0.45
48:EL:85:LEU:HA	48:EL:88:LYS:HG2	1.98	0.45
56:AK:68:ILE:HG13	56:AK:69:PRO:HD2	1.98	0.45
56:AK:154:LYS:HE3	56:AK:154:LYS:HB2	1.77	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
82:B0:40:LYS:HA	82:B0:40:LYS:HD2	1.83	0.45
4:BA:143:A:H2'	4:BA:144:A:C8	2.52	0.45
4:BA:232:U:H2'	4:BA:233:A:C8	2.52	0.45
4:BA:567:C:O2'	4:BA:1020:C:O2'	2.30	0.45
10:BJ:188:ARG:HE	10:BJ:192:ILE:HD11	1.82	0.45
10:BJ:220:SER:O	10:BJ:224:HIS:N	2.46	0.45
13:BN:160:GLN:O	13:BN:164:GLU:HG3	2.16	0.45
27:Bc:200:THR:O	27:Bc:204:ARG:HG3	2.17	0.45
35:Bk:189:HIS:CE1	35:Bk:194:PHE:HB2	2.50	0.45
35:Bk:203:GLU:O	35:Bk:206:GLU:HG3	2.17	0.45
41:Bq:115:ARG:HH21	41:Bq:119:ARG:NH2	2.15	0.45
56:AK:83:GLU:N	56:AK:83:GLU:OE2	2.50	0.45
58:AN:56:SER:O	58:AN:60:ASN:ND2	2.49	0.45
71:Ae:191:PHE:CA	71:Ae:196:ASN:HD22	2.29	0.45
80:Ao:423:MET:HG3	80:Ao:438:VAL:HG11	1.99	0.45
80:Ao:627:PRO:HB3	80:Ao:659:THR:HG21	1.99	0.45
4:BA:1006:U:H2'	4:BA:1007:G:O4'	2.17	0.44
4:BA:1388:A:H8	4:BA:1388:A:OP2	2.00	0.44
8:BF:83:HIS:HB2	8:BF:270:GLU:OE2	2.17	0.44
11:BK:25:ARG:HE	11:BK:65:PRO:HB3	1.83	0.44
15:BP:100:ARG:O	15:BP:104:LEU:HG	2.17	0.44
23:BX:140:ARG:HG3	27:Bc:77:VAL:HG21	2.00	0.44
25:Ba:236:LEU:HD23	25:Ba:341:VAL:HG13	1.99	0.44
25:Ba:283:TYR:HE2	25:Ba:285:TYR:HB2	1.82	0.44
27:Bc:137:LYS:O	27:Bc:141:ARG:HG3	2.17	0.44
27:Bc:219:GLU:HG2	27:Bc:222:VAL:HG12	1.99	0.44
44:Bv:128:MET:CE	88:B6:54:VAL:HG13	2.47	0.44
47:AA:243:G:N2	47:AA:256:G:OP1	2.44	0.44
47:AA:389:A:H5''	59:AO:101:GLN:HB2	1.99	0.44
50:AC:134:PHE:HD2	50:AC:135:LEU:HD22	1.82	0.44
51:AE:177:GLU:N	51:AE:177:GLU:OE2	2.50	0.44
52:AF:26:LEU:HD11	52:AF:39:LEU:HD22	1.98	0.44
52:AF:104:GLU:HG2	52:AF:104:GLU:O	2.18	0.44
54:AI:270:PHE:HD1	54:AI:287:TYR:HE1	1.65	0.44
48:FL:85:LEU:HA	48:FL:88:LYS:HG2	1.98	0.44
67:Aa:150:MET:HG3	67:Aa:151:PRO:HD2	1.99	0.44
71:Ae:250:LEU:HD21	76:Aj:145:HIS:CE1	2.53	0.44
71:Ae:261:LEU:HD11	71:Ae:302:VAL:HB	1.99	0.44
83:B1:11:TRP:O	83:B1:15:ARG:HG2	2.17	0.44
4:BA:176:U:H2'	4:BA:177:U:C6	2.53	0.44
4:BA:217:A:H62	36:Bl:111:ARG:HD2	1.83	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:BA:653:A:H2'	4:BA:654:G:O4'	2.17	0.44
16:BQ:39:PHE:HD2	16:BQ:166:ARG:HA	1.83	0.44
18:BS:96:HIS:HE2	18:BS:97:ARG:HH11	1.64	0.44
24:BY:84:ASN:HB3	24:BY:117:HIS:HD2	1.82	0.44
28:Bd:85:ILE:HG22	28:Bd:87:ASP:H	1.82	0.44
32:Bh:239:GLY:O	32:Bh:242:VAL:HG22	2.17	0.44
34:Bj:260:LEU:HD21	34:Bj:264:LEU:HD12	1.99	0.44
48:CL:40:LEU:HD13	48:FL:71:ILE:HG22	1.99	0.44
52:AF:25:THR:O	52:AF:29:LEU:HD23	2.17	0.44
58:AN:33:LYS:HD3	58:AN:36:ARG:HH21	1.82	0.44
69:Ac:91:GLU:OE2	70:Ad:120:ARG:NH2	2.50	0.44
73:Ag:341:PRO:HB2	77:Ak:311:LEU:HD21	1.99	0.44
4:BA:347:U:O4'	4:BA:349:A:H5'	2.17	0.44
4:BA:551:A:OP1	10:BJ:128:ASN:ND2	2.50	0.44
5:BB:41:G:H5''	18:BS:155:SER:HB2	1.99	0.44
7:BE:333:TYR:CD2	7:BE:339:GLN:HG2	2.53	0.44
9:BI:85:ASP:OD2	9:BI:87:LYS:HG2	2.16	0.44
10:BJ:33:VAL:HB	16:BQ:69:LEU:HD13	1.98	0.44
16:BQ:72:ILE:HD11	16:BQ:96:TYR:CE1	2.53	0.44
28:Bd:117:LEU:HD23	34:Bj:73:LEU:HD13	2.00	0.44
31:Bg:35:ARG:O	31:Bg:44:ARG:NH1	2.50	0.44
47:AA:210:U:H2'	47:AA:211:A:C8	2.53	0.44
47:AA:687:A:C8	58:AN:109:ILE:HD13	2.52	0.44
49:AB:177:ASN:HD21	49:AB:180:LEU:HD22	1.83	0.44
57:AL:70:PRO:HB3	57:AL:117:ASP:HB3	1.99	0.44
59:AO:176:TYR:HB3	61:AQ:90:GLN:HG3	2.00	0.44
82:B0:76:HIS:O	82:B0:130:PHE:HA	2.17	0.44
4:BA:266:C:H5''	4:BA:267:A:H5'	1.98	0.44
5:BB:23:A:OP2	26:Bb:106:ARG:NH1	2.48	0.44
8:BF:51:CYS:SG	8:BF:81:GLU:HG2	2.57	0.44
19:BT:60:PRO:HB2	71:Ae:127:HIS:ND1	2.32	0.44
21:BV:162:ALA:HB2	21:BV:199:ILE:HG13	1.98	0.44
24:BY:170:TRP:O	29:Be:75:ARG:NH2	2.50	0.44
25:Ba:193:LEU:HD11	45:Bw:183:SER:HB2	1.98	0.44
27:Bc:289:LEU:HD23	27:Bc:289:LEU:H	1.82	0.44
38:Bn:93:ARG:HG2	38:Bn:97:MET:HE3	2.00	0.44
45:Bw:110:THR:HB	45:Bw:112:THR:HG22	1.99	0.44
47:AA:893:A:H2'	47:AA:894:U:O4'	2.17	0.44
50:AC:81:HIS:ND1	50:AC:84:GLU:OE1	2.51	0.44
56:AK:138:ALA:HB1	56:AK:172:LEU:HD13	1.98	0.44
67:Aa:137:THR:O	67:Aa:141:VAL:HG23	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
77:Ak:105:LEU:O	77:Ak:109:LYS:HG3	2.18	0.44
80:Ao:332:LYS:HE3	80:Ao:332:LYS:HB2	1.78	0.44
83:B1:44:ARG:NE	83:B1:44:ARG:HA	2.32	0.44
2:B8:154:GLN:HE21	2:B8:154:GLN:HB2	1.61	0.44
4:BA:96:U:OP1	44:Bv:33:ARG:NH2	2.44	0.44
4:BA:191:A:N6	4:BA:1029:G:OP1	2.51	0.44
4:BA:570:A:H2'	4:BA:571:A:C8	2.53	0.44
4:BA:1056:G:H2'	4:BA:1324:C:O2'	2.18	0.44
5:BB:63:C:H2'	5:BB:64:A:C8	2.50	0.44
7:BE:56:GLU:O	17:BR:154:ARG:NH2	2.51	0.44
8:BF:283:LEU:HD21	15:BP:189:LYS:HG2	1.98	0.44
9:BI:139:LEU:O	9:BI:142:GLU:HG3	2.17	0.44
17:BR:43:GLU:OE1	87:B5:127:TYR:OH	2.35	0.44
19:BT:57:PRO:HA	19:BT:58:PRO:HD3	1.86	0.44
25:Ba:165:HIS:CD2	25:Ba:179:VAL:HG21	2.52	0.44
26:Bb:196:THR:OG1	26:Bb:324:ASP:OD2	2.30	0.44
45:Bw:399:ILE:HG22	45:Bw:404:VAL:HA	1.98	0.44
47:AA:332:U:H2'	47:AA:333:A:H8	1.83	0.44
47:AA:761:U:H2'	47:AA:762:G:C8	2.52	0.44
47:AA:938:U:H2'	47:AA:939:A:H8	1.81	0.44
49:AB:194:PHE:CZ	49:AB:207:ALA:HB3	2.53	0.44
52:AF:29:LEU:HD13	52:AF:78:MET:SD	2.57	0.44
73:Ag:135:LEU:HG	93:Ag:500:GTP:C4	2.53	0.44
73:Ag:173:ASN:O	73:Ag:176:ARG:CZ	2.65	0.44
77:Ak:99:MET:HE2	77:Ak:99:MET:HB3	1.88	0.44
77:Ak:322:LEU:HB3	77:Ak:324:LEU:HD13	2.00	0.44
81:Ap:59:LEU:HD11	81:Ap:119:ASN:HB2	1.99	0.44
4:BA:170:U:OP1	22:BW:205:THR:OG1	2.35	0.44
4:BA:1143:U:H2'	4:BA:1144:U:C6	2.53	0.44
4:BA:1279:G:H5''	42:Bt:14:GLY:H	1.81	0.44
6:BD:142:LEU:HD21	6:BD:149:LYS:HD2	1.98	0.44
6:BD:265:ARG:HE	6:BD:271:PRO:HD3	1.81	0.44
12:BL:110:TRP:CD1	12:BL:110:TRP:H	2.36	0.44
12:BL:118:LYS:O	12:BL:122:THR:HG22	2.17	0.44
21:BV:116:ASP:O	21:BV:198:ARG:HA	2.16	0.44
24:BY:199:MET:HG3	24:BY:204:ILE:HB	2.00	0.44
27:Bc:278:SER:H	27:Bc:295:GLY:HA2	1.82	0.44
30:Bf:67:ILE:HG21	32:Bh:258:THR:HG23	1.98	0.44
36:Bl:139:GLN:HB2	42:Bt:82:PHE:HE2	1.82	0.44
47:AA:299:U:H2'	47:AA:300:G:C8	2.53	0.44
47:AA:936:C:H2'	47:AA:937:A:C8	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
60:AP:63:GLU:OE1	60:AP:63:GLU:N	2.51	0.44
62:AR:68:LEU:O	62:AR:104:LYS:HE3	2.18	0.44
75:Ai:30:SER:O	75:Ai:33:VAL:HG22	2.18	0.44
77:Ak:68:TRP:CH2	77:Ak:84:PRO:HD2	2.53	0.44
77:Ak:268:LEU:HD21	77:Ak:321:LEU:HD11	1.99	0.44
80:Ao:395:ILE:HA	80:Ao:398:ASP:OD2	2.18	0.44
82:B0:102:GLU:HB2	82:B0:130:PHE:CD1	2.52	0.44
87:B5:159:LEU:HG	87:B5:174:ILE:HD11	1.99	0.44
4:BA:175:C:H2'	4:BA:176:U:C6	2.52	0.44
4:BA:852:C:OP2	6:BD:296:TYR:OH	2.32	0.44
4:BA:1501:C:H2'	4:BA:1502:U:C6	2.53	0.44
21:BV:178:PHE:HD1	21:BV:185:GLN:HG2	1.83	0.44
33:Bi:168:CYS:O	33:Bi:172:MET:HG2	2.17	0.44
34:Bj:215:PHE:CE2	34:Bj:229:ALA:HB3	2.51	0.44
38:Bn:57:TYR:OH	44:Bv:28:ARG:O	2.25	0.44
38:Bn:79:TRP:CE2	38:Bn:94:LYS:HD2	2.53	0.44
40:Bp:17:ARG:HE	40:Bp:65:ASP:HB2	1.83	0.44
46:Bx:43:ASP:N	46:Bx:43:ASP:OD1	2.51	0.44
47:AA:55:G:H2'	47:AA:56:U:C6	2.52	0.44
47:AA:760:A:H2'	47:AA:761:U:C6	2.53	0.44
48:CL:35:ILE:HG22	48:FL:76:LEU:HD22	1.99	0.44
48:EL:72:ALA:HA	48:HL:76:LEU:HD23	1.99	0.44
52:AF:78:MET:O	52:AF:82:LEU:HG	2.18	0.44
54:AI:88:LEU:HB2	77:Ak:85:LEU:HD11	2.00	0.44
67:Aa:298:VAL:HG23	67:Aa:329:LEU:HD13	1.98	0.44
72:Af:162:THR:HG22	72:Af:164:LEU:H	1.81	0.44
78:Am:48:GLU:HA	78:Am:51:VAL:HG22	2.00	0.44
4:BA:188:C:O2'	4:BA:190:U:OP2	2.23	0.44
4:BA:248:A:OP1	4:BA:323:A:N6	2.51	0.44
14:BO:69:VAL:HG23	14:BO:89:HIS:CE1	2.53	0.44
18:BS:110:TRP:CE2	18:BS:114:LYS:HD3	2.53	0.44
23:BX:42:PHE:CE2	23:BX:57:LEU:HD11	2.53	0.44
25:Ba:203:CYS:HB3	45:Bw:173:GLN:OE1	2.18	0.44
30:Bf:62:HIS:ND1	30:Bf:62:HIS:O	2.50	0.44
34:Bj:162:ARG:HH21	34:Bj:171:TRP:HB2	1.81	0.44
40:Bp:90:PHE:O	40:Bp:94:ILE:HG12	2.18	0.44
47:AA:154:C:H2'	47:AA:155:A:C8	2.53	0.44
47:AA:812:U:H2'	47:AA:813:G:H8	1.83	0.44
54:AI:201:LEU:HD11	54:AI:245:PHE:HA	2.00	0.44
54:AI:241:PHE:HE1	54:AI:244:ARG:HH21	1.66	0.44
67:Aa:254:VAL:HG21	67:Aa:289:HIS:HE1	1.83	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
73:Ag:66:HIS:CD2	73:Ag:99:MET:HB2	2.53	0.44
78:Am:29:LEU:HB3	78:Am:113:ASN:ND2	2.32	0.44
80:Ao:257:LYS:HD3	80:Ao:259:ALA:HB3	1.99	0.44
81:Ap:59:LEU:HD13	81:Ap:121:LYS:HB2	2.00	0.44
2:B8:113:ARG:HD3	15:BP:78:ILE:O	2.18	0.44
4:BA:52:A:H8	4:BA:347:U:H5	1.66	0.44
4:BA:568:C:C2	13:BN:78:SER:HB2	2.53	0.44
5:BB:25:G:H5''	18:BS:86:THR:HG21	2.00	0.44
7:BE:100:ILE:HG22	7:BE:296:LEU:HD12	1.99	0.44
10:BJ:81:LEU:HD11	41:Bq:115:ARG:HG3	2.00	0.44
12:BL:153:LYS:O	12:BL:157:MET:HG3	2.17	0.44
25:Ba:201:ARG:HD3	25:Ba:230:LEU:HD21	1.99	0.44
38:Bn:103:PHE:HB2	38:Bn:106:ASP:OD2	2.17	0.44
54:AI:73:PHE:O	54:AI:77:GLN:HG2	2.17	0.44
55:AJ:81:LYS:HE2	55:AJ:140:TYR:HE2	1.83	0.44
84:B2:96:SER:O	84:B2:100:LEU:HG	2.17	0.44
4:BA:122:G:N2	4:BA:125:A:OP2	2.46	0.43
4:BA:199:A:H5''	38:Bn:32:PHE:CD1	2.53	0.43
4:BA:364:A:O2'	4:BA:385:U:OP1	2.35	0.43
4:BA:1236:C:H2'	4:BA:1237:A:C8	2.53	0.43
4:BA:1270:G:H2'	4:BA:1323:U:H5''	2.00	0.43
12:BL:71:ASP:HB2	12:BL:155:ARG:HH11	1.83	0.43
15:BP:238:ARG:NH2	43:Bu:66:LYS:HD3	2.33	0.43
15:BP:252:LEU:HB3	15:BP:255:MET:HE3	1.99	0.43
22:BW:77:GLN:H	22:BW:171:HIS:HD2	1.64	0.43
26:Bb:51:TYR:CD2	82:B0:122:GLN:HA	2.53	0.43
33:Bi:231:LEU:HD11	33:Bi:234:GLY:O	2.18	0.43
34:Bj:173:LEU:HD22	34:Bj:233:PHE:HE2	1.83	0.43
44:Bv:128:MET:HE3	88:B6:54:VAL:HG13	2.00	0.43
45:Bw:147:GLU:OE1	45:Bw:173:GLN:HG3	2.17	0.43
47:AA:265:U:H2'	47:AA:266:U:C6	2.53	0.43
47:AA:364:U:H2'	47:AA:365:A:H8	1.83	0.43
47:AA:892:A:H2	76:Aj:17:ARG:HH11	1.65	0.43
51:AE:286:GLU:HB3	51:AE:288:HIS:CD2	2.52	0.43
56:AK:90:ILE:HD11	56:AK:163:ALA:HB3	1.98	0.43
67:Aa:338:GLU:HG3	67:Aa:341:LYS:HE2	2.00	0.43
4:BA:548:A:H2'	4:BA:549:C:C6	2.53	0.43
4:BA:742:U:H2'	4:BA:743:U:C6	2.52	0.43
6:BD:199:SER:HB2	6:BD:203:ARG:HD3	2.00	0.43
13:BN:16:ARG:HG3	32:Bh:267:LEU:HD11	2.00	0.43
24:BY:184:GLU:OE1	24:BY:186:THR:OG1	2.36	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:Bb:57:TYR:OH	82:B0:104:TYR:O	2.28	0.43
27:Bc:197:HIS:O	27:Bc:201:LYS:HG2	2.18	0.43
72:Af:109:VAL:HG12	72:Af:110:GLU:N	2.33	0.43
76:Aj:87:TRP:HE3	81:Ap:219:LEU:HD12	1.83	0.43
6:BD:178:ARG:HG2	6:BD:179:GLU:N	2.31	0.43
7:BE:94:SER:HA	7:BE:182:THR:HG21	1.99	0.43
11:BK:165:ALA:HA	11:BK:168:GLN:HG2	2.00	0.43
12:BL:185:ARG:O	12:BL:189:ARG:HG3	2.19	0.43
15:BP:156:VAL:O	15:BP:176:THR:HA	2.18	0.43
28:Bd:143:GLN:NE2	34:Bj:209:PRO:O	2.48	0.43
44:Bv:117:ARG:HH21	44:Bv:120:LEU:HD23	1.82	0.43
47:AA:73:G:H1	47:AA:155:A:H1'	1.84	0.43
47:AA:300:G:HO2'	59:AO:155:ARG:HH21	1.66	0.43
47:AA:579:A:OP1	58:AN:96:ARG:NH1	2.48	0.43
47:AA:925:A:O2'	47:AA:946:A:H5'	2.18	0.43
48:DL:83:ASN:C	48:DL:83:ASN:HD22	2.26	0.43
49:AB:72:PHE:HB2	49:AB:74:LEU:CD2	2.48	0.43
49:AB:128:LEU:HD11	49:AB:251:LEU:HD23	1.99	0.43
54:AI:102:GLU:HB3	54:AI:106:ARG:NH2	2.34	0.43
56:AK:156:LEU:HD11	63:AU:36:ILE:HG23	2.00	0.43
70:Ad:40:GLY:HA2	76:Aj:49:ARG:NH2	2.34	0.43
71:Ae:147:TRP:H	71:Ae:147:TRP:CD1	2.36	0.43
73:Ag:122:ARG:HG3	73:Ag:305:ILE:HB	2.01	0.43
73:Ag:283:PRO:HB2	73:Ag:293:ARG:HH21	1.83	0.43
78:Am:98:SER:OG	78:Am:99:LEU:N	2.51	0.43
83:B1:81:GLY:O	83:B1:130:ARG:NH1	2.50	0.43
4:BA:15:A:H5'	38:Bn:79:TRP:HH2	1.81	0.43
4:BA:447:A:H61	16:BQ:140:MET:HE2	1.83	0.43
4:BA:1393:C:H2'	4:BA:1394:U:O4'	2.18	0.43
13:BN:67:PHE:HE2	13:BN:92:ALA:HB1	1.83	0.43
25:Ba:212:THR:HG21	25:Ba:273:VAL:HA	2.01	0.43
27:Bc:306:PHE:O	27:Bc:310:ASN:ND2	2.45	0.43
33:Bi:226:ASP:OD1	33:Bi:227:ARG:N	2.51	0.43
34:Bj:132:LYS:HA	34:Bj:132:LYS:HD2	1.86	0.43
38:Bn:79:TRP:CZ2	38:Bn:94:LYS:HD2	2.53	0.43
47:AA:647:A:N6	63:AU:80:ARG:O	2.44	0.43
49:AB:105:LEU:HD21	49:AB:224:THR:HG22	1.98	0.43
73:Ag:55:PRO:HA	73:Ag:58:HIS:CD2	2.53	0.43
73:Ag:126:TYR:O	73:Ag:344:VAL:HG22	2.18	0.43
74:Ah:274:PHE:CE2	74:Ah:278:LEU:HD11	2.54	0.43
74:Ah:358:VAL:O	74:Ah:362:VAL:HG23	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
76:Aj:163:SER:HB3	76:Aj:190:MET:HB3	2.00	0.43
80:Ao:457:ARG:O	80:Ao:460:TYR:HB3	2.18	0.43
80:Ao:611:ALA:HB3	80:Ao:644:VAL:HB	1.99	0.43
4:BA:666:G:O6	4:BA:774:A:N6	2.52	0.43
4:BA:860:G:OP1	6:BD:104:TYR:OH	2.35	0.43
4:BA:942:U:H2'	4:BA:943:C:C6	2.52	0.43
4:BA:1002:U:C2	4:BA:1003:C:C5	3.06	0.43
4:BA:1531:C:H3'	4:BA:1532:U:H5''	2.01	0.43
8:BF:104:LYS:HA	38:Bn:54:PHE:HZ	1.84	0.43
13:BN:109:TYR:OH	13:BN:119:ARG:NE	2.52	0.43
13:BN:174:GLU:N	13:BN:174:GLU:OE2	2.51	0.43
14:BO:37:ARG:HG2	14:BO:54:PRO:HB2	1.99	0.43
16:BQ:240:ARG:NH2	16:BQ:246:TYR:OH	2.51	0.43
25:Ba:242:GLN:O	25:Ba:246:GLU:HG2	2.19	0.43
25:Ba:350:ARG:NH1	25:Ba:384:GLN:O	2.51	0.43
28:Bd:133:ARG:HA	28:Bd:136:ILE:HG22	2.00	0.43
34:Bj:145:ASP:OD1	34:Bj:147:THR:OG1	2.31	0.43
35:Bk:135:LEU:HB3	35:Bk:144:MET:HG3	2.00	0.43
47:AA:232:G:H5''	51:AE:196:ASN:HD22	1.83	0.43
47:AA:269:C:H2'	47:AA:270:C:C6	2.53	0.43
47:AA:489:G:H2'	47:AA:490:A:H8	1.84	0.43
47:AA:883:C:H1'	47:AA:886:A:N6	2.33	0.43
48:CL:46:LEU:O	48:CL:50:LEU:HG	2.18	0.43
61:AQ:65:LEU:HD12	61:AQ:66:LEU:N	2.33	0.43
71:Ae:122:ARG:HD3	71:Ae:159:TYR:CB	2.48	0.43
72:Af:145:LEU:HD21	72:Af:148:LEU:HD23	2.01	0.43
80:Ao:112:SER:HA	80:Ao:115:ASN:OD1	2.18	0.43
81:Ap:215:ARG:HG3	81:Ap:218:ARG:NH2	2.34	0.43
83:B1:134:LEU:HB3	83:B1:144:TYR:CE1	2.53	0.43
84:B2:104:TRP:CE3	84:B2:144:LEU:HD22	2.51	0.43
1:B7:74:ARG:HA	1:B7:74:ARG:HD3	1.75	0.43
4:BA:157:A:H2'	4:BA:158:A:C8	2.54	0.43
4:BA:165:A:N3	20:BU:52:LYS:HD3	2.34	0.43
4:BA:204:G:H2'	15:BP:40:PRO:HG3	2.01	0.43
4:BA:343:C:H2'	4:BA:344:U:C6	2.53	0.43
4:BA:529:G:P	41:Bq:115:ARG:HH22	2.41	0.43
4:BA:1213:A:O2'	88:B6:51:LYS:HB3	2.19	0.43
4:BA:1315:U:H2'	4:BA:1316:U:C6	2.52	0.43
8:BF:113:LYS:HD2	8:BF:157:GLY:HA2	2.01	0.43
15:BP:47:ARG:HB2	21:BV:183:ASN:HD22	1.83	0.43
15:BP:246:ASP:OD2	15:BP:249:LYS:NZ	2.43	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
21:BV:100:PHE:HB2	21:BV:111:LYS:HE2	2.01	0.43
24:BY:19:TYR:OH	24:BY:31:ASP:OD2	2.37	0.43
27:Bc:151:ILE:HG21	27:Bc:180:VAL:HG21	2.00	0.43
33:Bi:143:ASP:O	33:Bi:147:GLU:HG2	2.19	0.43
38:Bn:104:ILE:HD11	84:B2:212:ARG:HD3	1.99	0.43
47:AA:238:C:H2'	47:AA:239:U:H6	1.84	0.43
51:AE:369:HIS:HA	51:AE:387:PRO:HD3	2.00	0.43
59:AO:165:LYS:NZ	59:AO:193:PRO:HD3	2.33	0.43
60:AP:56:PRO:HA	60:AP:64:LYS:HZ3	1.84	0.43
76:Aj:4:LYS:HA	76:Aj:4:LYS:HD3	1.87	0.43
4:BA:232:U:H2'	4:BA:233:A:H8	1.83	0.43
4:BA:342:C:H2'	4:BA:343:C:C6	2.54	0.43
4:BA:611:U:H2'	4:BA:612:A:H8	1.84	0.43
4:BA:1050:C:O5'	4:BA:1051:G:H5''	2.19	0.43
6:BD:214:CYS:SG	6:BD:247:ARG:HG2	2.58	0.43
7:BE:64:LEU:HD21	27:Bc:311:LYS:HG3	2.00	0.43
8:BF:60:ARG:NH2	8:BF:61:PRO:O	2.50	0.43
8:BF:162:TYR:OH	38:Bn:71:LYS:NZ	2.32	0.43
13:BN:62:THR:HB	13:BN:129:PRO:HA	2.01	0.43
17:BR:60:ILE:HD11	17:BR:93:LEU:HD23	2.00	0.43
25:Ba:212:THR:O	25:Ba:326:ARG:NH1	2.44	0.43
25:Ba:272:ASP:OD1	25:Ba:272:ASP:N	2.52	0.43
28:Bd:187:PRO:HA	28:Bd:188:PRO:HD3	1.90	0.43
30:Bf:51:LEU:HD12	30:Bf:57:THR:O	2.19	0.43
47:AA:59:C:H3'	47:AA:60:C:C6	2.53	0.43
47:AA:189:A:H2'	47:AA:190:G:C8	2.52	0.43
47:AA:857:C:H2'	47:AA:858:C:C6	2.53	0.43
50:AC:76:LEU:O	58:AN:103:ARG:NH2	2.50	0.43
56:AK:165:LYS:HE2	56:AK:165:LYS:HB2	1.76	0.43
57:AL:129:LYS:HG3	57:AL:130:TYR:H	1.83	0.43
59:AO:218:VAL:O	59:AO:221:LEU:HG	2.19	0.43
64:AV:56:C:H2'	64:AV:57:C:C6	2.54	0.43
67:Aa:184:SER:HB2	67:Aa:187:ILE:HG12	2.00	0.43
70:Ad:94:LYS:HA	70:Ad:97:GLU:HG3	2.00	0.43
76:Aj:113:LYS:HA	76:Aj:113:LYS:HD3	1.76	0.43
4:BA:233:A:H2'	4:BA:234:G:C8	2.54	0.43
4:BA:316:U:H2'	4:BA:317:A:C8	2.54	0.43
4:BA:551:A:H5'	10:BJ:150:HIS:ND1	2.34	0.43
4:BA:1557:C:H4'	19:BT:86:ARG:HH22	1.84	0.43
7:BE:236:THR:HG22	7:BE:236:THR:O	2.18	0.43
11:BK:124:LYS:HE3	11:BK:124:LYS:HB3	1.90	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
16:BQ:96:TYR:HE1	16:BQ:153:PRO:HB3	1.83	0.43
47:AA:97:C:H2'	47:AA:98:A:O4'	2.18	0.43
47:AA:729:A:H2'	47:AA:730:C:O4'	2.19	0.43
48:CL:31:LEU:O	48:CL:35:ILE:HG13	2.18	0.43
49:AB:177:ASN:HB3	49:AB:181:LEU:HD23	1.99	0.43
51:AE:178:GLU:HG2	51:AE:181:ARG:HH21	1.84	0.43
52:AF:119:SER:OG	52:AF:120:THR:N	2.52	0.43
54:AI:87:HIS:O	54:AI:91:MET:HG3	2.18	0.43
69:Ac:50:PHE:O	69:Ac:53:PRO:HD2	2.18	0.43
71:Ae:323:LEU:HA	71:Ae:326:VAL:HG22	2.00	0.43
74:Ah:292:GLU:OE1	80:Ao:87:TYR:OH	2.32	0.43
76:Aj:109:LEU:O	76:Aj:109:LEU:HD23	2.18	0.43
80:Ao:301:ILE:HG13	80:Ao:302:GLU:N	2.33	0.43
80:Ao:353:LYS:NZ	80:Ao:392:SER:OG	2.52	0.43
81:Ap:211:ARG:HE	81:Ap:215:ARG:HH21	1.65	0.43
84:B2:130:SER:OG	84:B2:130:SER:O	2.36	0.43
4:BA:515:A:N1	4:BA:540:A:O2'	2.51	0.43
4:BA:516:G:H1'	4:BA:533:A:H2'	2.01	0.43
4:BA:698:U:H2'	4:BA:699:G:C8	2.53	0.43
4:BA:749:C:OP1	29:Be:47:LYS:NZ	2.44	0.43
4:BA:978:G:H4'	4:BA:1429:C:H4'	2.01	0.43
8:BF:200:PRO:HB2	8:BF:237:LEU:HD21	2.01	0.43
9:BI:100:GLN:HB2	9:BI:128:LEU:HD13	2.00	0.43
26:Bb:198:ALA:O	26:Bb:254:TYR:OH	2.30	0.43
37:Bm:114:PRO:HG2	37:Bm:117:ARG:HG3	2.00	0.43
39:Bo:65:ARG:O	39:Bo:69:GLN:HG3	2.18	0.43
43:Bu:117:LYS:HA	43:Bu:120:VAL:HG12	2.01	0.43
45:Bw:299:LYS:HA	45:Bw:299:LYS:HD3	1.85	0.43
47:AA:123:U:H2'	47:AA:124:A:C4	2.53	0.43
47:AA:704:U:H2'	47:AA:705:G:C8	2.53	0.43
47:AA:740:U:OP1	53:AG:197:ARG:NH2	2.46	0.43
47:AA:816:U:H2'	47:AA:817:G:O4'	2.19	0.43
51:AE:198:TRP:CE3	51:AE:201:VAL:HG21	2.53	0.43
51:AE:278:TYR:HB3	68:Ab:6:LEU:HD21	2.01	0.43
54:AI:350:ALA:CB	54:AI:370:LEU:HB3	2.49	0.43
64:AV:24:U:H2'	64:AV:25:C:H6	1.82	0.43
73:Ag:110:TYR:HB3	73:Ag:123:TYR:OH	2.18	0.43
83:B1:202:GLU:HG3	83:B1:203:TRP:CD1	2.53	0.43
85:B3:85:ILE:HG23	85:B3:109:LYS:HB3	2.01	0.43
4:BA:744:A:OP2	25:Ba:85:PRO:HB3	2.19	0.43
14:BO:107:LEU:HD23	14:BO:107:LEU:HA	1.90	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
15:BP:16:LEU:HD23	15:BP:19:LEU:HD12	2.00	0.43
17:BR:86:ILE:HB	17:BR:87:PRO:HD3	2.01	0.43
17:BR:156:ASP:OD2	27:Bc:315:ARG:NE	2.52	0.43
32:Bh:119:LEU:HD23	32:Bh:119:LEU:HA	1.89	0.43
36:Bl:150:LYS:NZ	42:Bt:79:ALA:O	2.49	0.43
37:Bm:138:ARG:HB2	37:Bm:142:ASP:HB2	1.99	0.43
43:Bu:185:GLU:O	43:Bu:189:LYS:HG2	2.18	0.43
44:Bv:80:GLU:HG2	44:Bv:81:GLN:N	2.33	0.43
44:Bv:120:LEU:HD12	44:Bv:120:LEU:HA	1.89	0.43
47:AA:72:A:H2'	47:AA:73:G:C8	2.54	0.43
47:AA:641:A:N7	51:AE:260:LYS:NZ	2.64	0.43
54:AI:125:MET:HE2	77:Ak:86:PRO:O	2.19	0.43
55:AJ:109:HIS:NE2	55:AJ:142:CYS:SG	2.88	0.43
68:Ab:119:VAL:O	68:Ab:123:LYS:HG2	2.19	0.43
71:Ae:169:LEU:HD13	71:Ae:177:ILE:HD11	2.01	0.43
80:Ao:520:LYS:O	80:Ao:527:ARG:NH2	2.51	0.43
88:B6:45:HIS:ND1	88:B6:46:TYR:N	2.67	0.43
4:BA:1059:C:H2'	4:BA:1060:C:C6	2.54	0.42
4:BA:1380:A:H5''	12:BL:102:ARG:NH2	2.32	0.42
6:BD:67:LYS:HB2	25:Ba:35:VAL:HG21	2.01	0.42
6:BD:181:ASP:HB3	6:BD:183:HIS:CD2	2.54	0.42
13:BN:25:MET:O	13:BN:149:ARG:NH1	2.52	0.42
19:BT:188:LEU:HD12	19:BT:191:ARG:NH1	2.34	0.42
26:Bb:329:ARG:HG3	26:Bb:329:ARG:NH1	2.34	0.42
27:Bc:153:ARG:HH22	27:Bc:256:ASP:C	2.22	0.42
32:Bh:159:PHE:O	32:Bh:162:SER:OG	2.29	0.42
33:Bi:160:LEU:O	33:Bi:164:VAL:HG22	2.19	0.42
34:Bj:187:THR:HG23	34:Bj:190:ARG:HH21	1.83	0.42
44:Bv:153:ARG:HE	44:Bv:153:ARG:HB2	1.71	0.42
52:AF:40:GLU:HG2	52:AF:42:LEU:HD22	2.01	0.42
58:AN:65:LYS:O	58:AN:69:GLU:HG2	2.18	0.42
60:AP:31:PHE:HE1	60:AP:53:SER:HB2	1.84	0.42
64:AV:3:U:H2'	64:AV:4:A:C8	2.54	0.42
71:Ae:119:ILE:HB	71:Ae:155:GLN:NE2	2.30	0.42
71:Ae:433:ARG:O	71:Ae:436:GLU:HG3	2.19	0.42
74:Ah:324:ILE:HG13	77:Ak:223:TYR:CE1	2.54	0.42
77:Ak:70:SER:HA	80:Ao:78:VAL:HG23	2.00	0.42
78:Am:48:GLU:OE1	78:Am:76:CYS:HB2	2.18	0.42
80:Ao:420:GLN:CG	80:Ao:456:ARG:HD2	2.41	0.42
80:Ao:586:PHE:CD2	80:Ao:591:ARG:HB3	2.50	0.42
87:B5:126:CYS:HA	87:B5:129:LYS:HE2	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:BA:193:U:H2'	4:BA:194:G:C8	2.54	0.42
4:BA:639:U:H5'	22:BW:147:ARG:NH1	2.34	0.42
4:BA:1017:U:O2'	4:BA:1440:U:H5'	2.18	0.42
7:BE:199:ARG:HD3	7:BE:199:ARG:HA	1.72	0.42
19:BT:84:ARG:HH11	19:BT:84:ARG:HG2	1.84	0.42
22:BW:188:LYS:HD2	22:BW:193:HIS:CE1	2.54	0.42
32:Bh:246:LYS:HE3	32:Bh:246:LYS:HB2	1.82	0.42
33:Bi:217:HIS:ND1	33:Bi:243:LEU:HD13	2.34	0.42
45:Bw:131:LEU:HD21	45:Bw:416:VAL:HG11	2.00	0.42
47:AA:59:C:H3'	47:AA:60:C:H6	1.85	0.42
47:AA:589:C:H2'	47:AA:590:C:C6	2.55	0.42
50:AC:87:VAL:HG21	75:Ai:53:PRO:HG2	2.00	0.42
53:AG:209:LEU:HD23	53:AG:209:LEU:HA	1.71	0.42
71:Ae:106:PRO:HB3	71:Ae:139:PRO:HB2	2.01	0.42
71:Ae:110:LEU:HD12	71:Ae:430:LEU:HB2	2.00	0.42
71:Ae:218:THR:HG23	71:Ae:251:PRO:O	2.19	0.42
74:Ah:294:ILE:HG13	80:Ao:71:LEU:HD23	2.01	0.42
77:Ak:87:VAL:HG13	77:Ak:107:LEU:HD11	2.01	0.42
80:Ao:398:ASP:OD1	80:Ao:399:ILE:N	2.53	0.42
80:Ao:536:MET:HE2	80:Ao:536:MET:HB3	1.92	0.42
4:BA:218:A:H4'	4:BA:219:A:H5'	2.01	0.42
4:BA:343:C:H2'	4:BA:344:U:H6	1.84	0.42
4:BA:523:U:O2'	4:BA:532:A:OP1	2.37	0.42
19:BT:114:GLY:HA3	19:BT:182:ARG:HG3	2.01	0.42
22:BW:188:LYS:HE3	22:BW:188:LYS:HB2	1.79	0.42
24:BY:209:ARG:HD3	24:BY:209:ARG:HA	1.89	0.42
25:Ba:201:ARG:NH1	25:Ba:418:TYR:O	2.38	0.42
28:Bd:164:ARG:NH2	35:Bk:164:ALA:HA	2.35	0.42
34:Bj:139:GLU:OE2	34:Bj:139:GLU:N	2.51	0.42
47:AA:285:C:N4	57:AL:43:LYS:O	2.48	0.42
47:AA:356:A:H2'	47:AA:357:G:H8	1.83	0.42
47:AA:902:C:H5''	79:An:171:ARG:NH1	2.35	0.42
47:AA:936:C:H2'	47:AA:937:A:H8	1.84	0.42
53:AG:163:LYS:HE2	53:AG:163:LYS:HB2	1.70	0.42
56:AK:66:PRO:HG3	63:AU:11:THR:HB	2.01	0.42
57:AL:71:LYS:HB3	57:AL:73:PRO:HD2	2.01	0.42
60:AP:29:ARG:NH1	60:AP:57:LEU:HB2	2.34	0.42
63:AU:72:ILE:HD12	78:Am:104:LEU:HD22	2.02	0.42
69:Ac:77:ARG:HG2	69:Ac:79:TYR:CZ	2.54	0.42
73:Ag:352:PHE:O	73:Ag:356:ILE:HG12	2.19	0.42
74:Ah:296:TRP:CB	74:Ah:302:LEU:HD23	2.49	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
74:Ah:329:TYR:CE1	74:Ah:362:VAL:HG22	2.54	0.42
77:Ak:192:LEU:O	77:Ak:196:VAL:HG22	2.18	0.42
82:B0:107:ASN:ND2	82:B0:108:PRO:HD2	2.34	0.42
85:B3:51:GLU:CD	85:B3:51:GLU:H	2.27	0.42
4:BA:441:A:H5''	16:BQ:66:ARG:HB2	2.01	0.42
4:BA:919:G:H2'	4:BA:920:C:C6	2.55	0.42
7:BE:142:MET:HE3	7:BE:180:ASN:HB3	2.01	0.42
8:BF:69:LEU:HD22	8:BF:193:LEU:HD12	2.01	0.42
8:BF:216:VAL:HG22	8:BF:258:THR:HB	2.00	0.42
10:BJ:176:LEU:HB3	10:BJ:188:ARG:HG3	2.02	0.42
11:BK:117:VAL:HG12	11:BK:141:VAL:HG23	2.01	0.42
11:BK:120:ILE:HA	11:BK:123:VAL:HG22	2.02	0.42
18:BS:96:HIS:CD2	18:BS:97:ARG:H	2.37	0.42
27:Bc:209:LYS:HE2	27:Bc:211:LEU:HD11	2.01	0.42
32:Bh:229:PHE:CZ	32:Bh:301:LEU:HD23	2.54	0.42
35:Bk:88:TYR:OH	35:Bk:162:LEU:HB3	2.19	0.42
43:Bu:142:GLN:O	43:Bu:146:LEU:HD23	2.19	0.42
45:Bw:143:CYS:O	45:Bw:147:GLU:HG2	2.19	0.42
45:Bw:265:LEU:N	45:Bw:266:PRO:HA	2.35	0.42
47:AA:82:C:H4'	61:AQ:5:ARG:HH21	1.84	0.42
47:AA:791:G:O2'	47:AA:817:G:O6	2.36	0.42
47:AA:875:A:N1	47:AA:893:A:C6	2.87	0.42
47:AA:882:A:O2'	71:Ae:140:ASN:OD1	2.37	0.42
53:AG:132:GLU:OE1	53:AG:132:GLU:N	2.49	0.42
54:AI:241:PHE:O	54:AI:244:ARG:HG2	2.19	0.42
67:Aa:355:LYS:HD3	67:Aa:365:GLU:OE1	2.19	0.42
69:Ac:27:VAL:HG22	69:Ac:106:LEU:HD12	2.00	0.42
73:Ag:87:ALA:O	73:Ag:91:LYS:HG2	2.19	0.42
73:Ag:125:LEU:HD23	73:Ag:342:ILE:HB	2.01	0.42
74:Ah:341:HIS:O	74:Ah:345:LEU:HD13	2.20	0.42
77:Ak:55:LEU:HB3	80:Ao:93:PRO:HB3	2.01	0.42
78:Am:11:ALA:HB2	78:Am:19:PRO:HB3	2.02	0.42
79:An:142:LYS:HD3	79:An:142:LYS:HA	1.82	0.42
80:Ao:109:ALA:O	80:Ao:112:SER:OG	2.31	0.42
85:B3:73:LYS:HE2	85:B3:73:LYS:HB3	1.91	0.42
13:BN:155:LEU:HD13	46:Bx:86:VAL:HG12	2.01	0.42
19:BT:152:ARG:HH12	19:BT:191:ARG:HG2	1.84	0.42
25:Ba:409:GLU:OE1	25:Ba:412:ARG:NH1	2.52	0.42
27:Bc:79:ASN:N	27:Bc:83:SER:OG	2.51	0.42
28:Bd:164:ARG:HH22	35:Bk:167:ALA:CB	2.32	0.42
30:Bf:109:ILE:HG12	30:Bf:123:TRP:HB2	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
47:AA:602:A:H2'	47:AA:603:U:C6	2.55	0.42
47:AA:689:G:O2'	75:Ai:75:HIS:NE2	2.42	0.42
47:AA:879:A:H2'	47:AA:880:U:C6	2.55	0.42
55:AJ:72:LEU:HB3	55:AJ:177:LEU:HD12	2.01	0.42
55:AJ:179:GLN:HG2	55:AJ:180:LEU:H	1.84	0.42
57:AL:129:LYS:HD3	57:AL:129:LYS:HA	1.80	0.42
60:AP:112:ARG:HB3	60:AP:116:ARG:NH2	2.35	0.42
67:Aa:250:LEU:HD23	67:Aa:289:HIS:HB3	2.00	0.42
69:Ac:136:LEU:C	69:Ac:137:ARG:HD3	2.45	0.42
71:Ae:192:ILE:HG13	71:Ae:224:TYR:OH	2.19	0.42
80:Ao:269:MET:CG	80:Ao:277:GLN:HE21	2.32	0.42
80:Ao:321:ASP:HA	80:Ao:324:LYS:HG2	2.01	0.42
48:HL:83:ASN:HD22	48:HL:83:ASN:C	2.26	0.42
83:B1:115:TYR:HD1	83:B1:122:ARG:HG2	1.84	0.42
88:B6:22:SER:OG	88:B6:54:VAL:HG11	2.19	0.42
4:BA:12:C:O2	4:BA:111:G:O2'	2.35	0.42
4:BA:363:G:H2'	4:BA:364:A:C8	2.55	0.42
4:BA:527:U:H2'	4:BA:529:G:N7	2.35	0.42
4:BA:687:U:O2'	4:BA:688:A:H5'	2.20	0.42
8:BF:141:ILE:HG13	8:BF:142:ARG:HD3	2.02	0.42
8:BF:286:PHE:HD2	38:Bn:66:PHE:CE2	2.37	0.42
15:BP:233:ARG:HE	15:BP:233:ARG:HB3	1.72	0.42
22:BW:130:HIS:HB2	22:BW:132:VAL:HG23	2.01	0.42
34:Bj:74:LEU:HD23	34:Bj:77:ILE:HD11	2.01	0.42
35:Bk:171:LEU:O	35:Bk:175:HIS:CD2	2.73	0.42
36:Bl:128:LEU:HB2	36:Bl:136:PRO:HG3	2.01	0.42
40:Bp:94:ILE:HG13	40:Bp:95:GLN:N	2.34	0.42
46:Bx:99:MET:HE3	46:Bx:100:LEU:N	2.31	0.42
47:AA:124:A:H2'	47:AA:125:U:C6	2.54	0.42
47:AA:388:U:H2'	47:AA:389:A:H8	1.83	0.42
47:AA:411:U:H2'	47:AA:412:A:H8	1.84	0.42
60:AP:67:ALA:HB2	67:Aa:218:TYR:OH	2.18	0.42
67:Aa:195:VAL:HG11	76:Aj:169:LEU:HD13	2.01	0.42
68:Ab:116:LYS:HD2	68:Ab:116:LYS:HA	1.83	0.42
71:Ae:370:GLN:HB3	71:Ae:374:ARG:HH22	1.84	0.42
76:Aj:85:TRP:CD1	76:Aj:85:TRP:N	2.88	0.42
76:Aj:180:LYS:HD2	76:Aj:180:LYS:HA	1.72	0.42
77:Ak:223:TYR:CZ	77:Ak:227:LEU:HD21	2.54	0.42
80:Ao:137:ILE:HG21	80:Ao:140:LEU:HD12	2.01	0.42
84:B2:87:TRP:HB3	84:B2:92:LEU:CD2	2.49	0.42
88:B6:41:LEU:HD12	88:B6:41:LEU:HA	1.91	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B8:175:ASP:HB3	2:B8:178:GLN:HB2	2.02	0.42
4:BA:269:U:H5''	25:Ba:395:LYS:HD3	2.01	0.42
4:BA:341:G:H2'	4:BA:342:C:C6	2.54	0.42
6:BD:129:ALA:HB1	6:BD:131:ARG:NH1	2.34	0.42
11:BK:112:VAL:HG23	11:BK:155:VAL:HG22	2.00	0.42
12:BL:191:ARG:HE	47:AA:244:C:H41	1.68	0.42
13:BN:170:TRP:NE1	46:Bx:92:PHE:HB3	2.34	0.42
14:BO:58:ILE:HD11	14:BO:76:ALA:HB2	2.01	0.42
28:Bd:146:ALA:HA	34:Bj:212:HIS:HB2	2.01	0.42
47:AA:378:A:H2'	47:AA:379:A:C8	2.54	0.42
56:AK:65:TYR:CE2	62:AR:131:LEU:HD11	2.55	0.42
56:AK:93:SER:OG	56:AK:94:TYR:N	2.52	0.42
62:AR:66:CYS:CB	62:AR:101:CYS:H	2.32	0.42
80:Ao:500:LEU:HD13	80:Ao:533:GLU:O	2.19	0.42
4:BA:489:C:C6	42:Bt:26:PRO:HG3	2.54	0.42
4:BA:695:U:H3	4:BA:696:A:N6	2.16	0.42
4:BA:869:G:H2'	4:BA:870:C:C6	2.55	0.42
7:BE:286:ASN:OD1	7:BE:287:GLY:N	2.53	0.42
11:BK:90:PHE:HD2	11:BK:120:ILE:HG23	1.84	0.42
11:BK:165:ALA:O	11:BK:168:GLN:HG2	2.20	0.42
15:BP:220:ARG:HA	15:BP:220:ARG:HD3	1.96	0.42
18:BS:165:MET:HE2	18:BS:165:MET:HB2	1.97	0.42
19:BT:189:TYR:CG	19:BT:243:ILE:HD11	2.55	0.42
25:Ba:381:LEU:HD23	25:Ba:381:LEU:HA	1.90	0.42
27:Bc:296:LEU:HD21	27:Bc:312:LEU:HB3	2.02	0.42
32:Bh:226:LYS:HE3	32:Bh:226:LYS:HB3	1.79	0.42
45:Bw:87:GLN:HE22	45:Bw:274:ASN:H	1.68	0.42
45:Bw:352:GLN:HB3	45:Bw:418:PHE:CD2	2.55	0.42
52:AF:119:SER:HB3	52:AF:121:LYS:HE3	2.01	0.42
62:AR:55:MET:HE1	62:AR:58:PRO:HG3	2.02	0.42
67:Aa:247:VAL:HG22	67:Aa:284:LEU:HD23	2.01	0.42
70:Ad:63:TYR:O	70:Ad:67:VAL:HG13	2.20	0.42
76:Aj:68:LEU:HD11	76:Aj:74:PHE:C	2.45	0.42
84:B2:209:ARG:HG3	84:B2:209:ARG:HH11	1.85	0.42
4:BA:140:A:H4'	24:BY:96:ARG:HB3	2.01	0.42
7:BE:96:ARG:HH11	7:BE:197:HIS:CD2	2.38	0.42
7:BE:175:LYS:HE2	7:BE:175:LYS:HB2	1.75	0.42
13:BN:28:PRO:HG3	13:BN:67:PHE:HE1	1.85	0.42
14:BO:75:LEU:HD23	14:BO:77:ILE:HD11	2.02	0.42
15:BP:51:ARG:NH1	15:BP:57:ARG:O	2.53	0.42
17:BR:134:LEU:HD23	87:B5:130:VAL:HG11	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
19:BT:95:GLU:O	19:BT:99:MET:HG2	2.20	0.42
20:BU:39:TYR:O	20:BU:43:VAL:HG23	2.20	0.42
23:BX:131:GLN:HA	23:BX:134:LYS:HG2	2.02	0.42
25:Ba:119:GLN:HG2	25:Ba:256:PHE:CZ	2.55	0.42
28:Bd:169:PHE:HB2	28:Bd:170:PRO:HD3	2.01	0.42
34:Bj:277:LEU:HD11	35:Bk:172:GLU:HG2	2.02	0.42
39:Bo:80:LEU:O	39:Bo:84:MET:HG3	2.20	0.42
47:AA:364:U:H2'	47:AA:365:A:C8	2.55	0.42
47:AA:791:G:H22	47:AA:818:A:P	2.42	0.42
47:AA:913:G:H2'	47:AA:914:A:C8	2.55	0.42
49:AB:77:LEU:HD21	49:AB:258:ARG:HH22	1.85	0.42
49:AB:252:PHE:O	49:AB:256:ILE:HG12	2.20	0.42
52:AF:5:GLU:OE2	62:AR:76:LYS:NZ	2.49	0.42
52:AF:40:GLU:OE1	70:Ad:187:TYR:OH	2.32	0.42
53:AG:53:LYS:HG3	53:AG:53:LYS:O	2.19	0.42
67:Aa:254:VAL:HA	67:Aa:264:TYR:HE1	1.85	0.42
69:Ac:51:ASN:HA	69:Ac:54:GLN:OE1	2.19	0.42
71:Ae:70:LEU:HD12	71:Ae:71:SER:N	2.35	0.42
71:Ae:137:HIS:C	76:Aj:104:TYR:HB2	2.45	0.42
76:Aj:174:ILE:O	76:Aj:178:ARG:HG2	2.20	0.42
79:An:151:ARG:O	79:An:155:ARG:HG3	2.20	0.42
80:Ao:579:LEU:HD11	80:Ao:599:LEU:HD12	2.00	0.42
80:Ao:598:MET:HE3	80:Ao:598:MET:HB3	1.94	0.42
88:B6:35:SER:OG	88:B6:38:ARG:HB2	2.20	0.42
4:BA:76:G:O2'	4:BA:88:G:O6	2.37	0.42
4:BA:354:U:H2'	4:BA:355:G:H8	1.84	0.42
4:BA:935:C:H2'	4:BA:936:A:C8	2.55	0.42
4:BA:1194:U:H3'	26:Bb:364:ARG:HE	1.84	0.42
7:BE:227:GLN:OE1	7:BE:227:GLN:N	2.53	0.42
17:BR:67:ASP:OD1	45:Bw:44:TYR:OH	2.31	0.42
18:BS:104:SER:O	18:BS:135:ARG:NH1	2.51	0.42
20:BU:29:ARG:HA	20:BU:29:ARG:HD3	1.88	0.42
21:BV:76:GLU:OE1	21:BV:80:HIS:NE2	2.53	0.42
23:BX:5:VAL:HG21	29:Be:135:ARG:HD2	2.01	0.42
24:BY:23:MET:HE2	24:BY:114:PRO:HB3	2.02	0.42
28:Bd:121:TRP:CD2	34:Bj:70:MET:HG3	2.55	0.42
35:Bk:122:GLU:N	35:Bk:158:GLN:O	2.53	0.42
40:Bp:40:GLU:HA	40:Bp:40:GLU:OE1	2.20	0.42
42:Bt:16:GLN:HB3	42:Bt:17:TRP:CD1	2.54	0.42
45:Bw:352:GLN:NE2	45:Bw:421:ASN:HD22	2.18	0.42
47:AA:483:U:H2'	47:AA:484:C:C6	2.55	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
47:AA:638:A:H2	51:AE:258:ILE:HD12	1.85	0.42
47:AA:858:C:H2'	47:AA:859:U:H6	1.84	0.42
50:AC:157:THR:OG1	80:Ao:92:ASP:OD1	2.33	0.42
53:AG:174:LEU:HD22	53:AG:178:ARG:HG2	2.01	0.42
55:AJ:74:LYS:CD	55:AJ:75:ARG:HG3	2.42	0.42
58:AN:66:ASP:HB3	74:Ah:378:ILE:HD11	2.01	0.42
69:Ac:139:CYS:O	69:Ac:142:GLU:HG2	2.20	0.42
80:Ao:161:ILE:HG13	80:Ao:162:LYS:N	2.35	0.42
80:Ao:481:TYR:CD1	80:Ao:485:ILE:HG13	2.54	0.42
3:B9:69:LEU:HB2	4:BA:560:A:N6	2.35	0.41
4:BA:470:C:H5'	10:BJ:43:GLU:OE2	2.19	0.41
4:BA:657:U:H1'	4:BA:658:U:C5	2.55	0.41
4:BA:1163:A:N6	4:BA:1164:G:O6	2.53	0.41
4:BA:1195:A:N6	4:BA:1208:A:O2'	2.52	0.41
4:BA:1528:A:H2'	4:BA:1529:A:H8	1.84	0.41
5:BB:50:A:H2'	5:BB:51:G:C8	2.55	0.41
7:BE:99:LEU:HD13	7:BE:198:PHE:CE2	2.55	0.41
10:BJ:81:LEU:O	10:BJ:84:ARG:HG2	2.20	0.41
10:BJ:181:ILE:O	10:BJ:184:THR:OG1	2.37	0.41
11:BK:138:SER:O	11:BK:141:VAL:HG12	2.20	0.41
15:BP:178:PHE:HB2	15:BP:214:TYR:CE2	2.55	0.41
20:BU:51:VAL:O	20:BU:55:ARG:HG3	2.20	0.41
30:Bf:46:LYS:O	30:Bf:62:HIS:HA	2.20	0.41
43:Bu:78:ILE:HD12	43:Bu:100:VAL:HG22	2.02	0.41
47:AA:388:U:H2'	47:AA:389:A:C8	2.55	0.41
47:AA:835:C:H2'	47:AA:836:G:H8	1.85	0.41
55:AJ:56:LYS:H	80:Ao:443:ASN:ND2	2.17	0.41
67:Aa:89:LYS:HE2	67:Aa:89:LYS:HB2	1.80	0.41
67:Aa:129:THR:HG22	67:Aa:130:GLN:N	2.34	0.41
74:Ah:298:LYS:HE3	74:Ah:298:LYS:HB3	1.66	0.41
81:Ap:90:THR:H	81:Ap:144:MET:HE2	1.85	0.41
85:B3:66:LEU:HD12	85:B3:122:LEU:HD23	2.02	0.41
85:B3:110:LEU:HD13	85:B3:119:ILE:HD13	2.02	0.41
2:B8:154:GLN:O	2:B8:158:LEU:HG	2.20	0.41
2:B8:170:ASN:HB3	2:B8:172:TYR:CE1	2.56	0.41
4:BA:219:A:P	15:BP:109:ARG:HH12	2.40	0.41
4:BA:230:G:H2'	4:BA:231:C:H6	1.84	0.41
4:BA:429:G:O2'	4:BA:599:A:N3	2.44	0.41
4:BA:480:G:OP2	21:BV:108:HIS:ND1	2.53	0.41
4:BA:906:G:H2'	4:BA:907:U:O4'	2.19	0.41
4:BA:1060:C:H2'	4:BA:1061:U:C6	2.54	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:BA:1177:U:H5''	16:BQ:144:LYS:HB2	2.01	0.41
5:BB:40:U:H2'	5:BB:41:G:H8	1.86	0.41
7:BE:198:PHE:HB3	7:BE:202:GLN:HE22	1.85	0.41
8:BF:184:GLN:NE2	15:BP:20:PRO:O	2.53	0.41
8:BF:290:TYR:HB2	8:BF:293:PHE:HB3	2.01	0.41
11:BK:124:LYS:HD3	11:BK:144:ILE:HD11	2.02	0.41
15:BP:43:ARG:HD2	21:BV:185:GLN:H	1.84	0.41
28:Bd:173:ARG:HB2	35:Bk:184:LEU:HD22	2.02	0.41
35:Bk:98:TYR:CE1	35:Bk:150:LEU:HD13	2.55	0.41
46:Bx:60:SER:HB2	46:Bx:75:TRP:CH2	2.56	0.41
47:AA:368:A:H2'	47:AA:368:A:N3	2.35	0.41
47:AA:441:U:H2'	47:AA:442:A:C8	2.54	0.41
47:AA:900:A:N3	47:AA:902:C:N4	2.67	0.41
49:AB:152:HIS:HA	49:AB:155:GLU:OE2	2.20	0.41
50:AC:118:GLU:OE2	50:AC:118:GLU:N	2.53	0.41
57:AL:138:LYS:HD3	57:AL:138:LYS:HA	1.86	0.41
64:AV:67:U:H2'	64:AV:68:A:C8	2.55	0.41
70:Ad:27:ARG:HB3	70:Ad:28:LYS:H	1.60	0.41
73:Ag:137:LEU:HD11	73:Ag:260:ALA:HB1	2.01	0.41
74:Ah:332:ASP:HB3	74:Ah:369:ARG:NH1	2.29	0.41
77:Ak:83:VAL:HG12	77:Ak:85:LEU:H	1.85	0.41
81:Ap:51:TYR:CE2	81:Ap:58:TYR:HB2	2.54	0.41
4:BA:344:U:H2'	4:BA:345:G:H8	1.85	0.41
4:BA:472:U:OP1	85:B3:73:LYS:HA	2.19	0.41
4:BA:1354:A:H61	4:BA:1465:A:N6	2.15	0.41
4:BA:1493:A:C6	19:BT:84:ARG:NH1	2.88	0.41
4:BA:1544:A:C2	4:BA:1545:A:C8	3.08	0.41
10:BJ:62:PRO:HA	10:BJ:65:LEU:HB3	2.02	0.41
15:BP:103:TYR:OH	36:Bl:93:ASN:ND2	2.48	0.41
16:BQ:195:LEU:O	16:BQ:199:ARG:HG3	2.20	0.41
17:BR:94:ALA:HB3	17:BR:95:PRO:HD3	2.03	0.41
20:BU:137:GLU:HB2	20:BU:138:PRO:HD3	2.02	0.41
25:Ba:210:SER:OG	25:Ba:223:HIS:ND1	2.37	0.41
25:Ba:233:LYS:HA	25:Ba:286:PRO:HD2	2.02	0.41
43:Bu:80:TYR:CE1	43:Bu:146:LEU:HD21	2.50	0.41
63:Au:66:MET:SD	78:Am:4:PRO:HD3	2.60	0.41
63:Au:82:ASP:N	63:Au:82:ASP:OD1	2.53	0.41
69:Ac:47:PHE:HA	69:Ac:51:ASN:HD22	1.86	0.41
71:Ae:203:VAL:HG12	71:Ae:207:ILE:HD11	2.02	0.41
48:GL:76:LEU:HD12	48:GL:76:LEU:HA	1.88	0.41
86:B4:55:VAL:HG22	86:B4:67:ILE:O	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:BA:196:U:O2'	4:BA:1014:A:OP1	2.36	0.41
4:BA:585:A:H5''	20:BU:58:ARG:NH1	2.35	0.41
4:BA:792:G:H2'	4:BA:793:A:H8	1.85	0.41
4:BA:802:G:N2	4:BA:987:G:OP2	2.40	0.41
19:BT:192:ASP:OD2	19:BT:229:TRP:HB3	2.21	0.41
26:Bb:138:GLU:HG2	26:Bb:141:ARG:NH1	2.35	0.41
32:Bh:166:VAL:HG11	32:Bh:192:ARG:HG3	2.02	0.41
34:Bj:200:MET:SD	34:Bj:241:GLY:HA3	2.60	0.41
44:Bv:30:PRO:HA	44:Bv:31:PRO:HD3	1.92	0.41
46:Bx:92:PHE:CZ	46:Bx:115:ILE:HG21	2.56	0.41
47:AA:633:C:OP1	51:AE:269:ARG:NH1	2.52	0.41
51:AE:136:ARG:N	75:Ai:67:PHE:O	2.53	0.41
67:Aa:259:PRO:HA	67:Aa:264:TYR:CD2	2.55	0.41
73:Ag:158:HIS:O	73:Ag:162:LYS:HG2	2.20	0.41
80:Ao:145:PHE:HA	80:Ao:163:LEU:HD13	2.02	0.41
80:Ao:514:GLN:HA	80:Ao:517:LYS:HZ3	1.86	0.41
4:BA:314:U:H2'	4:BA:315:G:H8	1.85	0.41
4:BA:421:U:H2'	4:BA:422:U:C6	2.55	0.41
4:BA:508:G:H4'	11:BK:151:LEU:HD12	2.03	0.41
4:BA:1121:U:H2'	4:BA:1122:A:H8	1.85	0.41
8:BF:249:ASN:ND2	8:BF:252:SER:OG	2.51	0.41
24:BY:18:ASN:HB3	24:BY:40:ARG:HB2	2.01	0.41
24:BY:161:ARG:CZ	24:BY:163:ASP:HB2	2.51	0.41
25:Ba:194:LYS:HB3	25:Ba:194:LYS:HE2	1.73	0.41
33:Bi:122:ILE:O	33:Bi:126:ARG:NE	2.53	0.41
38:Bn:105:ASP:OD1	38:Bn:106:ASP:N	2.52	0.41
43:Bu:162:GLN:OE1	43:Bu:163:PRO:HD2	2.21	0.41
47:AA:15:C:H2'	47:AA:16:U:C6	2.55	0.41
47:AA:470:A:OP1	51:AE:355:ARG:NH2	2.53	0.41
47:AA:572:A:OP2	47:AA:573:A:O2'	2.25	0.41
63:AU:73:ASN:HA	63:AU:76:MET:HG2	2.02	0.41
64:AV:3:U:H2'	64:AV:4:A:H8	1.85	0.41
67:Aa:231:ILE:HD12	67:Aa:236:ASN:HB3	2.02	0.41
67:Aa:270:HIS:O	67:Aa:273:GLU:HG2	2.21	0.41
71:Ae:290:LEU:HD12	71:Ae:290:LEU:HA	1.90	0.41
76:Aj:143:PRO:CB	76:Aj:145:HIS:CE1	3.03	0.41
77:Ak:107:LEU:HD23	77:Ak:107:LEU:HA	1.85	0.41
80:Ao:81:ASP:HB3	80:Ao:84:ALA:HB3	2.01	0.41
4:BA:1002:U:N3	4:BA:1003:C:C5	2.88	0.41
4:BA:1050:C:H41	4:BA:1323:U:H1'	1.85	0.41
4:BA:1344:C:O2'	4:BA:1387:A:H1'	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:BA:1420:A:H8	4:BA:1420:A:OP2	2.03	0.41
4:BA:1442:U:O2	7:BE:257:MET:HE1	2.21	0.41
8:BF:187:LEU:HD13	8:BF:259:LEU:HD23	2.02	0.41
15:BP:51:ARG:HD3	38:Bn:126:LYS:HD2	2.03	0.41
21:BV:105:PHE:HE1	39:Bo:49:TRP:HB3	1.86	0.41
28:Bd:150:LEU:HD12	34:Bj:212:HIS:CE1	2.55	0.41
31:Bg:146:ARG:HD3	31:Bg:146:ARG:HA	1.79	0.41
39:Bo:96:GLN:O	39:Bo:100:GLN:HG2	2.20	0.41
43:Bu:40:GLN:HB3	43:Bu:47:LYS:NZ	2.31	0.41
44:Bv:61:PHE:CD2	44:Bv:75:LEU:HD11	2.56	0.41
45:Bw:365:CYS:O	45:Bw:388:TRP:N	2.44	0.41
47:AA:88:C:H4'	60:AP:82:HIS:CE1	2.56	0.41
47:AA:512:A:H2'	47:AA:513:A:H8	1.85	0.41
48:CL:27:LYS:NZ	48:DL:70:ASP:HB3	2.35	0.41
48:CL:53:THR:OG1	48:CL:54:LEU:HD12	2.21	0.41
53:AG:203:GLU:O	53:AG:207:GLN:HG3	2.21	0.41
55:AJ:78:VAL:HG13	55:AJ:170:MET:HG3	2.01	0.41
80:Ao:511:MET:O	80:Ao:515:ILE:HG12	2.21	0.41
80:Ao:531:LYS:HE2	80:Ao:562:TYR:CE2	2.49	0.41
80:Ao:619:ASP:HA	80:Ao:622:LYS:HG2	2.02	0.41
83:B1:144:TYR:CE1	83:B1:148:THR:HG21	2.55	0.41
84:B2:229:GLU:O	84:B2:233:GLU:HG2	2.21	0.41
4:BA:464:U:OP1	42:Bt:23:ARG:NH1	2.54	0.41
4:BA:1021:C:H2'	4:BA:1022:G:C8	2.56	0.41
4:BA:1209:A:O3'	18:BS:66:ARG:NH1	2.53	0.41
4:BA:1278:C:H2'	4:BA:1279:G:C8	2.51	0.41
4:BA:1366:A:N7	4:BA:1537:A:O2'	2.50	0.41
10:BJ:142:ASN:ND2	48:CL:13:ALA:O	2.54	0.41
11:BK:127:ASP:O	11:BK:131:ALA:HB2	2.21	0.41
13:BN:21:LEU:HB3	13:BN:59:ILE:HG12	2.02	0.41
15:BP:68:GLU:OE1	15:BP:71:GLN:NE2	2.54	0.41
17:BR:146:LEU:HD11	27:Bc:304:ALA:HB2	2.03	0.41
20:BU:114:LYS:HD2	30:Bf:43:TYR:CE2	2.56	0.41
26:Bb:44:ASP:HB3	26:Bb:47:ARG:HG2	2.02	0.41
26:Bb:179:VAL:HG22	26:Bb:197:GLU:HG3	2.03	0.41
32:Bh:59:ARG:HH22	32:Bh:182:ALA:HA	1.86	0.41
45:Bw:233:ASN:HB2	45:Bw:293:PHE:HB2	2.02	0.41
45:Bw:356:THR:HG22	45:Bw:357:ASP:O	2.21	0.41
45:Bw:369:ASN:ND2	45:Bw:384:LYS:O	2.44	0.41
47:AA:277:A:H2'	47:AA:278:A:C8	2.56	0.41
47:AA:295:A:N3	47:AA:487:C:H1'	2.35	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
47:AA:302:U:OP1	59:AO:163:GLN:NE2	2.48	0.41
47:AA:578:G:OP1	55:AJ:128:LYS:NZ	2.43	0.41
47:AA:787:A:H2'	47:AA:788:G:C8	2.56	0.41
47:AA:835:C:H2'	47:AA:836:G:C8	2.55	0.41
49:AB:96:CYS:SG	72:Af:160:ASP:HB2	2.61	0.41
50:AC:156:GLN:HE22	77:Ak:108:LEU:HD11	1.85	0.41
52:AF:50:LYS:HE3	52:AF:57:ARG:CZ	2.51	0.41
55:AJ:155:VAL:HG11	77:Ak:132:CYS:HB3	2.01	0.41
71:Ae:405:LEU:HB3	71:Ae:406:PRO:HD3	2.03	0.41
71:Ae:436:GLU:O	71:Ae:439:GLU:HG2	2.20	0.41
83:B1:98:SER:OG	83:B1:99:LYS:N	2.52	0.41
4:BA:548:A:H2'	4:BA:549:C:H6	1.86	0.41
4:BA:568:C:O2'	46:Bx:164:ASN:O	2.26	0.41
4:BA:790:A:N6	4:BA:1000:A:O2'	2.53	0.41
4:BA:880:C:O2'	4:BA:922:A:H1'	2.21	0.41
4:BA:947:A:C6	14:BO:58:ILE:HG23	2.55	0.41
4:BA:1027:G:OP1	4:BA:1277:G:O2'	2.38	0.41
7:BE:334:ASP:HB3	7:BE:337:VAL:HG23	2.02	0.41
10:BJ:95:MET:HE2	10:BJ:156:SER:HB3	2.03	0.41
10:BJ:107:GLU:OE1	10:BJ:107:GLU:N	2.49	0.41
14:BO:87:VAL:HG13	14:BO:123:ILE:HG12	2.03	0.41
15:BP:247:ILE:HD13	43:Bu:59:TRP:CD1	2.55	0.41
25:Ba:210:SER:HB2	25:Ba:276:ASP:HB3	2.02	0.41
26:Bb:233:VAL:HG13	26:Bb:298:PHE:HB2	2.03	0.41
27:Bc:210:ASP:OD2	27:Bc:268:ARG:NH1	2.54	0.41
28:Bd:73:ARG:O	28:Bd:77:LYS:HG3	2.21	0.41
36:Bl:60:PRO:HA	36:Bl:61:PRO:HD3	1.97	0.41
41:Bq:122:ARG:O	41:Bq:126:ILE:HG12	2.20	0.41
45:Bw:147:GLU:HG3	45:Bw:170:PHE:CA	2.47	0.41
47:AA:603:U:H2'	47:AA:605:C:H4'	2.02	0.41
51:AE:89:PHE:CE1	75:Ai:62:MET:HG3	2.55	0.41
54:AI:99:PHE:CE2	54:AI:104:ILE:HG12	2.55	0.41
59:AO:71:SER:HA	59:AO:95:SER:HB2	2.03	0.41
59:AO:210:LEU:HB2	79:An:189:TRP:CZ2	2.56	0.41
68:Ab:27:GLU:N	68:Ab:27:GLU:OE1	2.53	0.41
74:Ah:354:PRO:HA	77:Ak:215:PRO:HG3	2.02	0.41
76:Aj:102:PRO:HA	76:Aj:112:GLY:HA3	2.03	0.41
83:B1:183:ARG:O	83:B1:187:ILE:HG12	2.20	0.41
84:B2:153:ASP:HA	84:B2:156:ARG:HG2	2.03	0.41
84:B2:158:LEU:HD12	84:B2:158:LEU:HA	1.93	0.41
2:B8:104:ARG:NH2	2:B8:160:LYS:O	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:BA:446:C:H2'	4:BA:447:A:C8	2.55	0.41
4:BA:777:A:H62	45:Bw:85:LYS:HE2	1.86	0.41
4:BA:1059:C:H2'	4:BA:1060:C:H6	1.85	0.41
4:BA:1060:C:H5''	4:BA:1152:G:O2'	2.21	0.41
4:BA:1131:U:H2'	4:BA:1132:G:C8	2.55	0.41
6:BD:124:GLU:HG2	6:BD:165:LEU:HD13	2.02	0.41
10:BJ:223:GLN:O	10:BJ:226:PRO:HD2	2.20	0.41
13:BN:154:ARG:HG2	46:Bx:127:LEU:HA	2.02	0.41
15:BP:189:LYS:HB3	15:BP:192:PRO:HG2	2.03	0.41
18:BS:61:VAL:HG12	18:BS:178:ILE:HB	2.03	0.41
18:BS:117:TYR:CE1	26:Bb:114:ARG:HG2	2.56	0.41
19:BT:138:ILE:HG23	19:BT:187:LEU:O	2.21	0.41
21:BV:140:VAL:HB	21:BV:147:LEU:HB2	2.03	0.41
25:Ba:60:TYR:HB3	25:Ba:70:LEU:HD12	2.02	0.41
25:Ba:110:ARG:HE	25:Ba:110:ARG:HB3	1.74	0.41
25:Ba:132:LEU:HB2	25:Ba:137:LEU:HD21	2.02	0.41
28:Bd:158:HIS:O	28:Bd:162:THR:HG23	2.21	0.41
30:Bf:75:ILE:HG23	32:Bh:289:VAL:HG11	2.02	0.41
32:Bh:170:ALA:HB1	32:Bh:175:VAL:HB	2.02	0.41
32:Bh:216:ARG:HA	32:Bh:220:ILE:HG13	2.03	0.41
33:Bi:137:PHE:N	33:Bi:138:PRO:HD2	2.36	0.41
33:Bi:158:ASP:O	33:Bi:162:THR:HG23	2.21	0.41
34:Bj:46:ARG:O	34:Bj:230:LYS:HG2	2.21	0.41
34:Bj:53:LEU:HD23	34:Bj:254:TRP:HH2	1.86	0.41
34:Bj:136:ARG:HE	34:Bj:136:ARG:HB3	1.65	0.41
44:Bv:29:PRO:HA	44:Bv:30:PRO:HD3	1.95	0.41
45:Bw:89:MET:SD	45:Bw:269:LYS:HB3	2.61	0.41
45:Bw:117:GLY:N	45:Bw:389:GLY:O	2.50	0.41
47:AA:308:A:C2	69:Ac:36:THR:HG21	2.56	0.41
47:AA:669:C:O2'	53:AG:35:SER:O	2.33	0.41
49:AB:65:PRO:HA	49:AB:71:PHE:CD1	2.55	0.41
49:AB:79:SER:H	49:AB:82:SER:HG	1.66	0.41
51:AE:316:CYS:HB2	51:AE:321:MET:HG3	2.02	0.41
52:AF:39:LEU:HD13	52:AF:66:VAL:HG22	2.02	0.41
53:AG:37:TYR:CE2	54:AI:375:PRO:HD2	2.52	0.41
54:AI:238:GLU:O	54:AI:242:VAL:HG23	2.21	0.41
56:AK:183:ILE:HG23	56:AK:183:ILE:O	2.21	0.41
57:AL:104:GLU:O	57:AL:106:HIS:HD2	2.04	0.41
60:AP:54:TYR:CD1	60:AP:66:VAL:HG22	2.56	0.41
61:AQ:28:VAL:HG11	61:AQ:66:LEU:CD1	2.49	0.41
48:FL:76:LEU:HD12	48:FL:76:LEU:HA	1.88	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
62:AR:117:ILE:HD12	62:AR:117:ILE:HA	1.92	0.41
67:Aa:246:HIS:O	67:Aa:250:LEU:HD13	2.20	0.41
68:Ab:125:LEU:HD23	68:Ab:125:LEU:HA	1.93	0.41
69:Ac:78:PHE:CD2	69:Ac:106:LEU:HD11	2.56	0.41
71:Ae:124:ASP:HA	71:Ae:127:HIS:CD2	2.56	0.41
71:Ae:235:ASP:OD1	71:Ae:235:ASP:N	2.50	0.41
71:Ae:360:VAL:O	71:Ae:360:VAL:HG12	2.21	0.41
72:Af:153:ARG:NH1	72:Af:157:ALA:O	2.52	0.41
73:Ag:126:TYR:OH	73:Ag:343:LEU:HB2	2.21	0.41
73:Ag:162:LYS:HB3	73:Ag:315:LEU:HD11	2.03	0.41
75:Ai:66:ARG:HA	75:Ai:71:TYR:O	2.20	0.41
75:Ai:86:LYS:HG3	75:Ai:89:ARG:NH2	2.36	0.41
76:Aj:58:VAL:HA	76:Aj:137:HIS:HD1	1.85	0.41
80:Ao:400:MET:O	80:Ao:404:MET:HG2	2.21	0.41
80:Ao:442:LEU:HD12	80:Ao:443:ASN:N	2.35	0.41
80:Ao:617:PHE:CD2	80:Ao:633:LEU:HD11	2.55	0.41
81:Ap:136:HIS:CD2	81:Ap:139:TYR:HB2	2.56	0.41
83:B1:216:ARG:HG3	83:B1:216:ARG:HH11	1.85	0.41
84:B2:153:ASP:OD1	84:B2:156:ARG:NH2	2.54	0.41
1:B7:78:GLY:O	1:B7:82:ILE:HG12	2.21	0.41
4:BA:467:A:OP1	42:Bt:15:ARG:NH2	2.54	0.41
4:BA:725:C:H5	25:Ba:389:LEU:H	1.69	0.41
4:BA:749:C:N3	23:BX:51:VAL:HG12	2.36	0.41
8:BF:184:GLN:HE22	15:BP:22:VAL:HG23	1.86	0.41
10:BJ:217:GLN:O	10:BJ:221:MET:HG2	2.21	0.41
11:BK:115:LYS:HD2	41:Bq:92:TYR:CE2	2.56	0.41
12:BL:183:MET:HG3	47:AA:918:A:O2'	2.21	0.41
13:BN:32:ALA:HB1	13:BN:108:ILE:HG13	2.03	0.41
19:BT:116:ILE:HD12	19:BT:178:LYS:HB3	2.03	0.41
20:BU:125:GLU:HB3	21:BV:68:ILE:HD12	2.03	0.41
25:Ba:344:SER:OG	25:Ba:355:LEU:HB3	2.20	0.41
30:Bf:126:ARG:HD3	30:Bf:126:ARG:HA	1.89	0.41
31:Bg:55:ALA:HB2	31:Bg:63:ILE:HD12	2.03	0.41
32:Bh:228:LEU:HB2	32:Bh:307:PHE:CD2	2.56	0.41
47:AA:756:U:H2'	47:AA:757:U:C6	2.56	0.41
52:AF:49:TYR:CE1	52:AF:51:ILE:HG12	2.56	0.41
57:AL:104:GLU:OE1	57:AL:104:GLU:HA	2.21	0.41
61:AQ:94:PRO:HG3	69:Ac:77:ARG:HH21	1.85	0.41
71:Ae:406:PRO:O	71:Ae:409:GLU:HG2	2.21	0.41
77:Ak:250:MET:SD	77:Ak:254:ILE:HD11	2.60	0.41
80:Ao:666:LEU:HD12	80:Ao:666:LEU:HA	1.94	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
81:Ap:94:CYS:SG	81:Ap:95:ILE:HG12	2.61	0.41
48:HL:76:LEU:HD12	48:HL:76:LEU:HA	1.88	0.41
4:BA:41:A:H2'	4:BA:42:C:H4'	2.02	0.40
4:BA:446:C:O2'	4:BA:447:A:O5'	2.36	0.40
4:BA:871:A:H8	4:BA:871:A:OP2	2.04	0.40
4:BA:1165:U:OP2	4:BA:1166:A:O2'	2.33	0.40
5:BB:1:G:O6	5:BB:70:A:N6	2.54	0.40
5:BB:50:A:H2'	5:BB:51:G:H8	1.85	0.40
7:BE:155:PHE:HA	27:Bc:305:HIS:NE2	2.36	0.40
8:BF:220:ASP:O	8:BF:245:ALA:N	2.54	0.40
11:BK:54:ASN:O	11:BK:58:LYS:HB3	2.21	0.40
11:BK:113:THR:OG1	11:BK:114:LEU:N	2.54	0.40
12:BL:201:ARG:NH2	47:AA:231:G:N3	2.69	0.40
13:BN:11:TRP:HZ2	32:Bh:267:LEU:HB2	1.86	0.40
18:BS:143:PHE:CD1	18:BS:175:PRO:HD3	2.56	0.40
23:BX:10:TYR:HB3	29:Be:54:MET:HE2	2.03	0.40
25:Ba:156:ASN:ND2	45:Bw:285:PRO:HG3	2.36	0.40
28:Bd:72:ILE:HD13	28:Bd:72:ILE:HA	1.87	0.40
32:Bh:60:ARG:HD2	32:Bh:60:ARG:HA	1.83	0.40
47:AA:254:C:H41	57:AL:74:ASN:HD22	1.69	0.40
47:AA:702:G:H5'	58:AN:36:ARG:HD3	2.03	0.40
50:AC:138:TYR:CZ	50:AC:142:LEU:HD22	2.56	0.40
51:AE:300:ARG:HD3	51:AE:300:ARG:HA	1.82	0.40
60:AP:31:PHE:CE1	60:AP:53:SER:HB2	2.56	0.40
69:Ac:114:GLU:HA	69:Ac:117:GLU:HG2	2.03	0.40
73:Ag:122:ARG:HA	73:Ag:305:ILE:O	2.22	0.40
76:Aj:170:LEU:HA	76:Aj:173:MET:SD	2.61	0.40
80:Ao:254:MET:HE3	80:Ao:256:GLU:HB2	2.03	0.40
80:Ao:447:ASN:HB3	80:Ao:448:TRP:HE3	1.86	0.40
4:BA:269:U:H2'	4:BA:270:U:C6	2.56	0.40
4:BA:385:U:H2'	4:BA:386:U:H6	1.85	0.40
4:BA:658:U:H2'	4:BA:659:U:H6	1.86	0.40
4:BA:743:U:O3'	25:Ba:88:ARG:NH2	2.54	0.40
4:BA:1083:C:H2'	4:BA:1084:C:C6	2.56	0.40
8:BF:230:VAL:HG12	8:BF:242:LEU:HD22	2.03	0.40
9:BI:75:ARG:NH1	83:B1:100:ARG:NH2	2.67	0.40
11:BK:130:PHE:HA	11:BK:133:GLN:HG2	2.03	0.40
13:BN:67:PHE:HB2	13:BN:72:TRP:CD1	2.56	0.40
17:BR:149:LEU:HB3	27:Bc:312:LEU:HD22	2.04	0.40
19:BT:178:LYS:HG3	19:BT:207:ILE:HD12	2.03	0.40
23:BX:10:TYR:HA	23:BX:14:ASN:ND2	2.27	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
27:Bc:163:LEU:HB3	27:Bc:178:TYR:CE2	2.56	0.40
27:Bc:177:CYS:SG	27:Bc:294:GLN:HG2	2.61	0.40
33:Bi:92:GLU:OE1	33:Bi:92:GLU:N	2.54	0.40
47:AA:928:A:O2'	47:AA:929:A:H5'	2.21	0.40
49:AB:214:MET:HE3	49:AB:214:MET:HB3	1.94	0.40
51:AE:163:GLU:O	51:AE:167:LYS:HG2	2.20	0.40
51:AE:168:VAL:O	51:AE:172:MET:HG2	2.21	0.40
55:Aj:104:ILE:CG2	55:Aj:145:LEU:HD22	2.52	0.40
55:Aj:158:GLU:OE2	74:Ah:305:PHE:HB3	2.22	0.40
62:AR:82:SER:HB2	62:AR:126:LYS:HD3	2.03	0.40
68:Ab:34:ILE:HD12	68:Ab:34:ILE:H	1.86	0.40
71:Ae:167:TYR:CD1	71:Ae:170:LYS:HE2	2.56	0.40
73:Ag:267:LEU:HB2	73:Ag:268:TRP:CE3	2.55	0.40
73:Ag:345:SER:OG	73:Ag:346:ASN:N	2.54	0.40
73:Ag:364:TRP:NE1	73:Ag:394:CYS:O	2.54	0.40
77:Ak:133:THR:HG22	80:Ao:63:LYS:HE3	2.04	0.40
80:Ao:492:HIS:CG	80:Ao:493:SER:N	2.89	0.40
80:Ao:578:SER:HA	80:Ao:581:TYR:CD2	2.57	0.40
83:B1:141:PHE:CZ	83:B1:145:ILE:HD11	2.55	0.40
4:BA:476:U:H2'	4:BA:477:A:H8	1.85	0.40
4:BA:1125:U:H2'	4:BA:1126:U:C6	2.56	0.40
4:BA:1183:U:C6	4:BA:1230:U:H2'	2.56	0.40
8:BF:68:SER:O	8:BF:210:ARG:NH2	2.38	0.40
8:BF:249:ASN:OD1	8:BF:249:ASN:N	2.54	0.40
14:BO:41:VAL:HG12	14:BO:105:VAL:O	2.21	0.40
25:Ba:367:ASP:O	25:Ba:371:LYS:NZ	2.42	0.40
26:Bb:351:HIS:CE1	26:Bb:368:ARG:HB3	2.56	0.40
31:Bg:28:ARG:HD2	32:Bh:60:ARG:HH21	1.86	0.40
31:Bg:32:SER:OG	31:Bg:68:ARG:HD3	2.22	0.40
32:Bh:238:MET:O	32:Bh:242:VAL:HG13	2.21	0.40
34:Bj:235:LYS:HD2	34:Bj:275:PHE:CD1	2.56	0.40
37:Bm:89:ASP:OD1	37:Bm:125:ARG:NH2	2.54	0.40
47:AA:421:A:H3'	47:AA:422:C:H6	1.86	0.40
47:AA:578:G:H3'	47:AA:579:A:H8	1.86	0.40
47:AA:643:G:O2'	47:AA:651:G:OP2	2.25	0.40
50:AC:164:TYR:CE2	50:AC:165:LYS:HE3	2.57	0.40
51:AE:157:ILE:HD11	80:Ao:116:ALA:HA	2.04	0.40
52:AF:26:LEU:HD22	52:AF:36:VAL:HG21	2.03	0.40
52:AF:114:GLU:HA	52:AF:114:GLU:OE1	2.21	0.40
55:Aj:70:ASP:OD1	55:Aj:71:THR:N	2.53	0.40
55:Aj:92:GLU:HG3	55:Aj:108:VAL:HG11	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
64:AV:44:U:H4'	64:AV:45:G:H5''	2.02	0.40
67:Aa:309:ASP:OD1	67:Aa:313:ARG:HD2	2.22	0.40
70:Ad:197:VAL:HG21	70:Ad:199:ARG:HH21	1.86	0.40
71:Ae:119:ILE:HG23	71:Ae:124:ASP:HB2	2.04	0.40
71:Ae:204:VAL:HG22	71:Ae:221:LEU:HD22	2.02	0.40
73:Ag:267:LEU:HD21	73:Ag:292:LEU:HD23	2.04	0.40
80:Ao:512:VAL:HG11	80:Ao:537:LEU:HD11	2.03	0.40
80:Ao:602:PHE:CD2	80:Ao:609:PRO:HD3	2.55	0.40
4:BA:248:A:H2'	4:BA:249:G:C8	2.57	0.40
4:BA:396:C:H1'	4:BA:415:A:N6	2.36	0.40
4:BA:800:A:H2'	4:BA:801:G:O4'	2.22	0.40
4:BA:1083:C:H2'	4:BA:1084:C:H6	1.86	0.40
4:BA:1521:U:C5	13:BN:177:ARG:HB3	2.49	0.40
17:BR:41:ARG:NH2	87:B5:134:THR:OG1	2.54	0.40
21:BV:148:LEU:HB2	30:Bf:58:ILE:HB	2.02	0.40
35:Bk:110:VAL:HG13	35:Bk:157:VAL:HG21	2.04	0.40
38:Bn:99:GLY:O	38:Bn:102:MET:HG3	2.21	0.40
38:Bn:103:PHE:CE2	83:B1:2:PRO:HG2	2.57	0.40
47:AA:300:G:N7	92:AA:3001:SPM:N10	2.67	0.40
50:AC:41:ARG:NH1	50:AC:41:ARG:HB2	2.37	0.40
52:AF:100:GLN:N	52:AF:100:GLN:OE1	2.55	0.40
53:AG:131:ALA:HA	53:AG:134:GLN:OE1	2.22	0.40
57:AL:51:PRO:HA	57:AL:89:ARG:HH12	1.86	0.40
67:Aa:135:GLU:HA	67:Aa:138:LYS:HG2	2.04	0.40
71:Ae:101:TYR:CD2	71:Ae:134:LYS:HE3	2.57	0.40
71:Ae:142:TRP:HD1	76:Aj:74:PHE:HD2	1.68	0.40
71:Ae:322:VAL:O	71:Ae:325:ARG:HG2	2.22	0.40
72:Af:129:GLU:H	72:Af:129:GLU:CD	2.30	0.40
73:Ag:66:HIS:NE2	73:Ag:99:MET:HB2	2.36	0.40
74:Ah:328:LYS:H	74:Ah:328:LYS:HG2	1.71	0.40
75:Ai:6:GLU:O	75:Ai:10:ARG:HG2	2.21	0.40
76:Aj:170:LEU:O	76:Aj:173:MET:HB2	2.21	0.40
77:Ak:196:VAL:HB	77:Ak:199:ARG:HG3	2.03	0.40
80:Ao:163:LEU:HD23	80:Ao:163:LEU:HA	1.95	0.40
81:Ap:90:THR:OG1	81:Ap:144:MET:SD	2.79	0.40
4:BA:36:A:C4	29:Be:32:LYS:HG3	2.57	0.40
4:BA:573:A:C6	46:Bx:196:HIS:HB2	2.56	0.40
4:BA:668:A:H2'	4:BA:669:C:C6	2.57	0.40
4:BA:851:G:H1'	6:BD:207:TYR:HE1	1.86	0.40
4:BA:1502:U:H2'	4:BA:1503:C:C6	2.56	0.40
7:BE:206:VAL:HG13	7:BE:297:VAL:CG1	2.51	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:BF:102:TRP:O	8:BF:106:PHE:HD1	2.05	0.40
14:BO:77:ILE:HD12	14:BO:107:LEU:HD21	2.04	0.40
15:BP:203:ARG:HH21	15:BP:262:PRO:HA	1.86	0.40
15:BP:264:GLN:NE2	15:BP:269:LEU:O	2.55	0.40
25:Ba:203:CYS:H	45:Bw:173:GLN:HE22	1.70	0.40
44:Bv:86:GLU:O	44:Bv:90:ARG:HG2	2.21	0.40
44:Bv:111:GLU:O	44:Bv:115:GLN:HG2	2.21	0.40
47:AA:883:C:H3'	71:Ae:433:ARG:NH1	2.34	0.40
51:AE:180:ASP:O	51:AE:184:LYS:HG2	2.22	0.40
56:AK:177:ILE:HD11	63:AU:19:VAL:HG13	2.03	0.40
60:AP:38:HIS:CD2	60:AP:39:SER:N	2.90	0.40
61:AQ:57:GLN:O	61:AQ:87:LYS:NZ	2.46	0.40
48:FL:70:ASP:O	48:FL:74:LEU:HG	2.22	0.40
71:Ae:101:TYR:CE2	71:Ae:134:LYS:HG2	2.56	0.40
80:Ao:385:HIS:O	80:Ao:388:PRO:HD2	2.21	0.40
48:GL:70:ASP:O	48:GL:74:LEU:HG	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all 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	
1	B7	44/95 (46%)	44 (100%)	0	0	100	100
2	B8	93/188 (50%)	93 (100%)	0	0	100	100
3	B9	36/100 (36%)	36 (100%)	0	0	100	100
6	BD	238/306 (78%)	231 (97%)	7 (3%)	0	100	100
7	BE	305/399 (76%)	291 (95%)	14 (5%)	0	100	100
8	BF	248/294 (84%)	242 (98%)	6 (2%)	0	100	100
9	BI	96/268 (36%)	95 (99%)	1 (1%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
10	BJ	210/262 (80%)	199 (95%)	11 (5%)	0	100	100
11	BK	174/192 (91%)	172 (99%)	2 (1%)	0	100	100
12	BL	135/184 (73%)	133 (98%)	2 (2%)	0	100	100
13	BN	175/178 (98%)	171 (98%)	4 (2%)	0	100	100
14	BO	113/145 (78%)	110 (97%)	3 (3%)	0	100	100
15	BP	286/296 (97%)	276 (96%)	10 (4%)	0	100	100
16	BQ	220/251 (88%)	217 (99%)	3 (1%)	0	100	100
17	BR	151/169 (89%)	146 (97%)	5 (3%)	0	100	100
18	BS	141/180 (78%)	131 (93%)	10 (7%)	0	100	100
19	BT	238/292 (82%)	233 (98%)	4 (2%)	1 (0%)	30	59
20	BU	138/149 (93%)	138 (100%)	0	0	100	100
21	BV	153/209 (73%)	149 (97%)	4 (3%)	0	100	100
22	BW	164/210 (78%)	158 (96%)	6 (4%)	0	100	100
23	BX	147/150 (98%)	145 (99%)	2 (1%)	0	100	100
24	BY	204/216 (94%)	197 (97%)	7 (3%)	0	100	100
25	Ba	391/423 (92%)	378 (97%)	13 (3%)	0	100	100
26	Bb	352/380 (93%)	333 (95%)	19 (5%)	0	100	100
27	Bc	293/334 (88%)	281 (96%)	11 (4%)	1 (0%)	36	65
28	Bd	136/206 (66%)	129 (95%)	7 (5%)	0	100	100
29	Be	120/135 (89%)	112 (93%)	8 (7%)	0	100	100
30	Bf	106/142 (75%)	102 (96%)	2 (2%)	2 (2%)	6	26
31	Bg	146/159 (92%)	137 (94%)	9 (6%)	0	100	100
32	Bh	287/332 (86%)	274 (96%)	13 (4%)	0	100	100
33	Bi	258/306 (84%)	244 (95%)	14 (5%)	0	100	100
34	Bj	211/279 (76%)	200 (95%)	11 (5%)	0	100	100
35	Bk	151/269 (56%)	144 (95%)	7 (5%)	0	100	100
36	Bl	131/166 (79%)	128 (98%)	3 (2%)	0	100	100
37	Bm	107/198 (54%)	106 (99%)	1 (1%)	0	100	100
38	Bn	95/128 (74%)	90 (95%)	5 (5%)	0	100	100
39	Bo	95/124 (77%)	94 (99%)	1 (1%)	0	100	100
40	Bp	95/112 (85%)	90 (95%)	5 (5%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
41	Bq	52/138 (38%)	49 (94%)	3 (6%)	0	100	100
42	Bt	92/102 (90%)	86 (94%)	6 (6%)	0	100	100
43	Bu	147/205 (72%)	139 (95%)	7 (5%)	1 (1%)	18	47
44	Bv	133/222 (60%)	133 (100%)	0	0	100	100
45	Bw	385/433 (89%)	365 (95%)	20 (5%)	0	100	100
46	Bx	160/196 (82%)	152 (95%)	8 (5%)	0	100	100
48	CL	43/198 (22%)	43 (100%)	0	0	100	100
48	DL	25/198 (13%)	25 (100%)	0	0	100	100
48	EL	26/198 (13%)	26 (100%)	0	0	100	100
48	FL	25/198 (13%)	25 (100%)	0	0	100	100
48	GL	25/198 (13%)	25 (100%)	0	0	100	100
48	HL	24/198 (12%)	24 (100%)	0	0	100	100
49	AB	218/289 (75%)	211 (97%)	7 (3%)	0	100	100
50	AC	130/167 (78%)	123 (95%)	7 (5%)	0	100	100
51	AE	341/430 (79%)	329 (96%)	12 (4%)	0	100	100
52	AF	120/276 (44%)	119 (99%)	1 (1%)	0	100	100
53	AG	206/242 (85%)	204 (99%)	2 (1%)	0	100	100
54	AI	326/397 (82%)	320 (98%)	6 (2%)	0	100	100
55	AJ	138/200 (69%)	125 (91%)	13 (9%)	0	100	100
56	AK	135/196 (69%)	130 (96%)	5 (4%)	0	100	100
57	AL	107/139 (77%)	99 (92%)	8 (8%)	0	100	100
58	AN	99/128 (77%)	98 (99%)	1 (1%)	0	100	100
59	AO	173/239 (72%)	169 (98%)	4 (2%)	0	100	100
60	AP	115/135 (85%)	112 (97%)	3 (3%)	0	100	100
61	AQ	110/130 (85%)	108 (98%)	2 (2%)	0	100	100
62	AR	95/143 (66%)	95 (100%)	0	0	100	100
63	AU	84/87 (97%)	82 (98%)	2 (2%)	0	100	100
66	AZ	16/18 (89%)	15 (94%)	1 (6%)	0	100	100
67	Aa	290/382 (76%)	284 (98%)	6 (2%)	0	100	100
68	Ab	133/190 (70%)	131 (98%)	2 (2%)	0	100	100
69	Ac	167/173 (96%)	166 (99%)	1 (1%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
70	Ad	175/205 (85%)	175 (100%)	0	0	100	100
71	Ae	386/455 (85%)	361 (94%)	23 (6%)	2 (0%)	24	54
72	Af	97/188 (52%)	95 (98%)	2 (2%)	0	100	100
73	Ag	351/410 (86%)	337 (96%)	14 (4%)	0	100	100
74	Ah	118/387 (30%)	117 (99%)	1 (1%)	0	100	100
75	Ai	97/106 (92%)	95 (98%)	2 (2%)	0	100	100
76	Aj	211/218 (97%)	205 (97%)	6 (3%)	0	100	100
77	Ak	273/325 (84%)	266 (97%)	7 (3%)	0	100	100
78	Am	114/118 (97%)	109 (96%)	5 (4%)	0	100	100
79	An	70/199 (35%)	69 (99%)	1 (1%)	0	100	100
80	Ao	564/699 (81%)	548 (97%)	16 (3%)	0	100	100
81	Ap	188/258 (73%)	182 (97%)	6 (3%)	0	100	100
82	B0	108/148 (73%)	107 (99%)	1 (1%)	0	100	100
83	B1	242/256 (94%)	241 (100%)	1 (0%)	0	100	100
84	B2	177/252 (70%)	172 (97%)	5 (3%)	0	100	100
85	B3	116/161 (72%)	113 (97%)	3 (3%)	0	100	100
86	B4	60/126 (48%)	55 (92%)	5 (8%)	0	100	100
87	B5	108/188 (57%)	107 (99%)	1 (1%)	0	100	100
88	B6	50/65 (77%)	49 (98%)	1 (2%)	0	100	100
All	All	14328/19635 (73%)	13864 (97%)	457 (3%)	7 (0%)	100	100

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
19	BT	68	PRO
30	Bf	80	PRO
43	Bu	167	PRO
30	Bf	78	PRO
71	Ae	342	PRO
71	Ae	341	GLN
27	Bc	107	GLY

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all 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	
1	B7	41/78 (53%)	41 (100%)	0	100	100
2	B8	87/162 (54%)	86 (99%)	1 (1%)	65	74
3	B9	36/77 (47%)	36 (100%)	0	100	100
6	BD	193/248 (78%)	193 (100%)	0	100	100
7	BE	263/320 (82%)	263 (100%)	0	100	100
8	BF	217/251 (86%)	217 (100%)	0	100	100
9	BI	88/228 (39%)	88 (100%)	0	100	100
10	BJ	192/230 (84%)	192 (100%)	0	100	100
11	BK	129/151 (85%)	129 (100%)	0	100	100
12	BL	125/162 (77%)	125 (100%)	0	100	100
13	BN	156/157 (99%)	156 (100%)	0	100	100
14	BO	99/123 (80%)	99 (100%)	0	100	100
15	BP	245/249 (98%)	245 (100%)	0	100	100
16	BQ	190/210 (90%)	190 (100%)	0	100	100
17	BR	132/143 (92%)	132 (100%)	0	100	100
18	BS	123/153 (80%)	123 (100%)	0	100	100
19	BT	212/258 (82%)	212 (100%)	0	100	100
20	BU	118/127 (93%)	118 (100%)	0	100	100
21	BV	136/178 (76%)	136 (100%)	0	100	100
22	BW	144/180 (80%)	144 (100%)	0	100	100
23	BX	116/134 (87%)	116 (100%)	0	100	100
24	BY	185/192 (96%)	185 (100%)	0	100	100
25	Ba	348/365 (95%)	348 (100%)	0	100	100
26	Bb	310/328 (94%)	310 (100%)	0	100	100
27	Bc	271/299 (91%)	270 (100%)	1 (0%)	84	83
28	Bd	127/181 (70%)	127 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
29	Be	100/108 (93%)	100 (100%)	0	100	100
30	Bf	80/133 (60%)	80 (100%)	0	100	100
31	Bg	128/136 (94%)	128 (100%)	0	100	100
32	Bh	251/284 (88%)	251 (100%)	0	100	100
33	Bi	236/275 (86%)	236 (100%)	0	100	100
34	Bj	190/242 (78%)	190 (100%)	0	100	100
35	Bk	135/226 (60%)	135 (100%)	0	100	100
36	Bl	122/147 (83%)	122 (100%)	0	100	100
37	Bm	103/178 (58%)	103 (100%)	0	100	100
38	Bn	88/113 (78%)	88 (100%)	0	100	100
39	Bo	77/97 (79%)	77 (100%)	0	100	100
40	Bp	79/88 (90%)	79 (100%)	0	100	100
41	Bq	50/114 (44%)	50 (100%)	0	100	100
42	Bt	75/82 (92%)	75 (100%)	0	100	100
43	Bu	126/177 (71%)	126 (100%)	0	100	100
44	Bv	115/183 (63%)	115 (100%)	0	100	100
45	Bw	340/373 (91%)	340 (100%)	0	100	100
46	Bx	149/173 (86%)	149 (100%)	0	100	100
48	CL	30/157 (19%)	30 (100%)	0	100	100
48	DL	26/157 (17%)	26 (100%)	0	100	100
48	EL	27/157 (17%)	27 (100%)	0	100	100
48	FL	26/157 (17%)	26 (100%)	0	100	100
48	GL	26/157 (17%)	26 (100%)	0	100	100
48	HL	25/157 (16%)	25 (100%)	0	100	100
49	AB	187/233 (80%)	187 (100%)	0	100	100
50	AC	115/142 (81%)	115 (100%)	0	100	100
51	AE	282/351 (80%)	278 (99%)	4 (1%)	59	70
52	AF	107/210 (51%)	107 (100%)	0	100	100
53	AG	181/205 (88%)	181 (100%)	0	100	100
54	AI	273/333 (82%)	271 (99%)	2 (1%)	76	78
55	AJ	130/180 (72%)	130 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
56	AK	103/151 (68%)	103 (100%)	0	100	100
57	AL	92/116 (79%)	92 (100%)	0	100	100
58	AN	92/114 (81%)	92 (100%)	0	100	100
59	AO	159/205 (78%)	159 (100%)	0	100	100
60	AP	97/113 (86%)	97 (100%)	0	100	100
61	AQ	97/114 (85%)	97 (100%)	0	100	100
62	AR	89/127 (70%)	89 (100%)	0	100	100
63	AU	77/78 (99%)	77 (100%)	0	100	100
67	Aa	258/330 (78%)	258 (100%)	0	100	100
68	Ab	113/162 (70%)	113 (100%)	0	100	100
69	Ac	152/155 (98%)	152 (100%)	0	100	100
70	Ad	149/168 (89%)	149 (100%)	0	100	100
71	Ae	325/393 (83%)	325 (100%)	0	100	100
72	Af	86/160 (54%)	86 (100%)	0	100	100
73	Ag	312/361 (86%)	312 (100%)	0	100	100
74	Ah	109/346 (32%)	109 (100%)	0	100	100
75	Ai	86/93 (92%)	86 (100%)	0	100	100
76	Aj	188/190 (99%)	188 (100%)	0	100	100
77	Ak	249/289 (86%)	249 (100%)	0	100	100
78	Am	100/102 (98%)	100 (100%)	0	100	100
79	An	66/174 (38%)	66 (100%)	0	100	100
80	Ao	478/611 (78%)	477 (100%)	1 (0%)	87	85
81	Ap	170/225 (76%)	170 (100%)	0	100	100
82	B0	90/115 (78%)	90 (100%)	0	100	100
83	B1	219/229 (96%)	219 (100%)	0	100	100
84	B2	164/228 (72%)	164 (100%)	0	100	100
85	B3	110/150 (73%)	110 (100%)	0	100	100
86	B4	45/114 (40%)	45 (100%)	0	100	100
87	B5	99/163 (61%)	99 (100%)	0	100	100
88	B6	49/60 (82%)	49 (100%)	0	100	100
All	All	12585/16745 (75%)	12576 (100%)	9 (0%)	87	89

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	B8	154	GLN
27	Bc	170	PRO
51	AE	174	GLN
51	AE	187	VAL
51	AE	188	LYS
51	AE	190	GLU
54	AI	248	SER
54	AI	251	ILE
80	Ao	329	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (131) such sidechains are listed below:

Mol	Chain	Res	Type
2	B8	154	GLN
6	BD	166	ASN
6	BD	169	HIS
6	BD	206	GLN
6	BD	236	GLN
6	BD	253	HIS
7	BE	154	HIS
7	BE	180	ASN
8	BF	74	GLN
8	BF	83	HIS
8	BF	153	HIS
10	BJ	151	ASN
10	BJ	239	GLN
12	BL	125	ASN
12	BL	143	GLN
12	BL	182	ASN
13	BN	9	GLN
13	BN	89	GLN
14	BO	104	ASN
14	BO	142	GLN
14	BO	143	ASN
15	BP	102	GLN
15	BP	219	ASN
16	BQ	138	GLN
17	BR	69	ASN
17	BR	109	GLN
17	BR	112	ASN
18	BS	162	GLN

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Mol	Chain	Res	Type
19	BT	239	ASN
22	BW	73	HIS
22	BW	77	GLN
22	BW	199	GLN
23	BX	27	GLN
25	Ba	214	ASN
25	Ba	324	GLN
26	Bb	77	HIS
26	Bb	220	ASN
26	Bb	295	GLN
28	Bd	143	GLN
29	Be	50	GLN
29	Be	121	GLN
30	Bf	62	HIS
31	Bg	58	ASN
31	Bg	127	GLN
32	Bh	42	GLN
32	Bh	65	ASN
32	Bh	106	GLN
32	Bh	172	ASN
32	Bh	177	GLN
33	Bi	193	GLN
33	Bi	196	GLN
33	Bi	235	GLN
33	Bi	251	HIS
34	Bj	67	GLN
34	Bj	87	HIS
34	Bj	198	ASN
35	Bk	111	HIS
35	Bk	138	GLN
35	Bk	158	GLN
35	Bk	175	HIS
35	Bk	177	ASN
35	Bk	189	HIS
36	Bl	92	HIS
36	Bl	93	ASN
37	Bm	68	GLN
37	Bm	107	ASN
38	Bn	69	HIS
38	Bn	120	HIS
40	Bp	15	GLN
40	Bp	57	HIS

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Mol	Chain	Res	Type
44	Bv	79	GLN
45	Bw	186	ASN
45	Bw	352	GLN
49	AB	152	HIS
49	AB	177	ASN
50	AC	60	HIS
50	AC	115	ASN
50	AC	154	HIS
50	AC	156	GLN
54	AI	148	HIS
54	AI	156	GLN
54	AI	327	HIS
54	AI	385	GLN
55	AJ	130	HIS
55	AJ	147	HIS
57	AL	105	HIS
58	AN	55	ASN
59	AO	78	GLN
59	AO	201	HIS
60	AP	59	ASN
60	AP	106	ASN
62	AR	83	GLN
67	Aa	243	GLN
67	Aa	245	GLN
67	Aa	268	HIS
67	Aa	324	HIS
67	Aa	327	HIS
67	Aa	369	GLN
68	Ab	66	HIS
69	Ac	51	ASN
70	Ad	65	GLN
70	Ad	113	ASN
70	Ad	161	GLN
71	Ae	137	HIS
71	Ae	155	GLN
71	Ae	196	ASN
73	Ag	62	HIS
73	Ag	89	GLN
73	Ag	227	GLN
73	Ag	301	GLN
74	Ah	295	GLN
74	Ah	331	GLN

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Mol	Chain	Res	Type
74	Ah	387	ASN
76	Aj	88	GLN
77	Ak	277	ASN
78	Am	24	ASN
78	Am	71	GLN
78	Am	105	ASN
79	An	164	GLN
79	An	166	GLN
80	Ao	273	GLN
80	Ao	420	GLN
80	Ao	458	ASN
80	Ao	492	HIS
80	Ao	507	ASN
80	Ao	580	ASN
82	B0	72	HIS
82	B0	107	ASN
82	B0	110	ASN
85	B3	102	ASN
87	B5	115	HIS

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
4	BA	1542/1571 (98%)	410 (26%)	2 (0%)
47	AA	959/962 (99%)	178 (18%)	0
5	BB	64/73 (87%)	15 (23%)	0
64	AV	70/71 (98%)	22 (31%)	0
65	AX	5/6 (83%)	1 (20%)	1 (20%)
All	All	2640/2683 (98%)	626 (23%)	3 (0%)

All (626) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
4	BA	4	A
4	BA	7	G
4	BA	11	G
4	BA	19	U
4	BA	20	A
4	BA	21	C
4	BA	22	U
4	BA	26	C

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Mol	Chain	Res	Type
4	BA	27	A
4	BA	31	A
4	BA	32	C
4	BA	36	A
4	BA	37	A
4	BA	40	C
4	BA	41	A
4	BA	45	A
4	BA	46	A
4	BA	49	A
4	BA	56	A
4	BA	57	A
4	BA	59	A
4	BA	60	U
4	BA	63	A
4	BA	66	U
4	BA	67	A
4	BA	68	A
4	BA	82	G
4	BA	83	A
4	BA	96	U
4	BA	97	A
4	BA	104	C
4	BA	105	G
4	BA	109	U
4	BA	115	U
4	BA	118	U
4	BA	119	A
4	BA	122	G
4	BA	129	A
4	BA	130	A
4	BA	132	G
4	BA	139	G
4	BA	140	A
4	BA	141	A
4	BA	142	U
4	BA	143	A
4	BA	146	A
4	BA	147	U
4	BA	149	A
4	BA	163	C
4	BA	164	A

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Mol	Chain	Res	Type
4	BA	165	A
4	BA	168	A
4	BA	172	C
4	BA	178	C
4	BA	180	A
4	BA	182	C
4	BA	186	U
4	BA	190	U
4	BA	192	A
4	BA	205	A
4	BA	218	A
4	BA	219	A
4	BA	223	A
4	BA	225	C
4	BA	229	A
4	BA	231	C
4	BA	237	A
4	BA	238	C
4	BA	239	C
4	BA	243	A
4	BA	245	A
4	BA	254	G
4	BA	263	G
4	BA	271	U
4	BA	272	A
4	BA	273	A
4	BA	275	A
4	BA	277	A
4	BA	295	G
4	BA	309	A
4	BA	311	A
4	BA	312	C
4	BA	322	G
4	BA	323	A
4	BA	324	G
4	BA	329	A
4	BA	330	A
4	BA	331	A
4	BA	336	A
4	BA	337	A
4	BA	338	C
4	BA	340	A

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Mol	Chain	Res	Type
4	BA	351	A
4	BA	352	G
4	BA	359	G
4	BA	366	A
4	BA	368	A
4	BA	369	G
4	BA	373	U
4	BA	374	U
4	BA	376	A
4	BA	389	A
4	BA	390	A
4	BA	393	A
4	BA	394	C
4	BA	398	A
4	BA	403	C
4	BA	409	A
4	BA	414	A
4	BA	417	G
4	BA	427	G
4	BA	428	A
4	BA	433	G
4	BA	440	A
4	BA	446	C
4	BA	447	A
4	BA	448	G
4	BA	459	A
4	BA	460	C
4	BA	467	A
4	BA	468	A
4	BA	473	G
4	BA	474	A
4	BA	479	A
4	BA	480	G
4	BA	488	U
4	BA	490	U
4	BA	491	U
4	BA	492	A
4	BA	493	A
4	BA	494	U
4	BA	497	U
4	BA	498	A
4	BA	499	C

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Mol	Chain	Res	Type
4	BA	500	C
4	BA	501	A
4	BA	503	A
4	BA	505	U
4	BA	509	C
4	BA	514	A
4	BA	515	A
4	BA	516	G
4	BA	517	C
4	BA	518	A
4	BA	526	A
4	BA	527	U
4	BA	530	A
4	BA	532	A
4	BA	533	A
4	BA	541	A
4	BA	547	A
4	BA	548	A
4	BA	549	C
4	BA	550	A
4	BA	551	A
4	BA	553	U
4	BA	554	U
4	BA	557	C
4	BA	560	A
4	BA	561	C
4	BA	562	A
4	BA	563	U
4	BA	570	A
4	BA	574	A
4	BA	576	U
4	BA	578	A
4	BA	579	U
4	BA	584	A
4	BA	586	C
4	BA	592	G
4	BA	593	C
4	BA	595	C
4	BA	596	A
4	BA	617	A
4	BA	618	A
4	BA	625	A

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Mol	Chain	Res	Type
4	BA	626	G
4	BA	631	A
4	BA	633	U
4	BA	634	G
4	BA	640	A
4	BA	648	C
4	BA	649	A
4	BA	654	G
4	BA	665	C
4	BA	673	C
4	BA	674	U
4	BA	675	U
4	BA	683	U
4	BA	684	A
4	BA	689	A
4	BA	694	A
4	BA	695	U
4	BA	697	C
4	BA	704	U
4	BA	707	A
4	BA	713	A
4	BA	719	C
4	BA	720	C
4	BA	722	A
4	BA	723	A
4	BA	724	A
4	BA	725	C
4	BA	727	A
4	BA	728	C
4	BA	744	A
4	BA	745	A
4	BA	746	U
4	BA	748	A
4	BA	763	C
4	BA	764	A
4	BA	774	A
4	BA	777	A
4	BA	778	A
4	BA	780	G
4	BA	782	A
4	BA	783	A
4	BA	784	G

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Mol	Chain	Res	Type
4	BA	819	A
4	BA	825	C
4	BA	835	A
4	BA	838	A
4	BA	847	U
4	BA	848	C
4	BA	852	C
4	BA	854	U
4	BA	855	U
4	BA	859	A
4	BA	864	U
4	BA	872	A
4	BA	883	G
4	BA	887	C
4	BA	889	C
4	BA	892	G
4	BA	894	U
4	BA	895	U
4	BA	896	A
4	BA	897	A
4	BA	899	G
4	BA	900	G
4	BA	902	C
4	BA	909	U
4	BA	910	U
4	BA	911	C
4	BA	922	A
4	BA	923	A
4	BA	924	G
4	BA	925	G
4	BA	926	U
4	BA	931	U
4	BA	932	A
4	BA	933	A
4	BA	935	C
4	BA	938	U
4	BA	939	U
4	BA	948	A
4	BA	950	U
4	BA	958	U
4	BA	959	G
4	BA	960	U

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Mol	Chain	Res	Type
4	BA	961	A
4	BA	962	U
4	BA	964	A
4	BA	965	A
4	BA	966	U
4	BA	967	G
4	BA	970	C
4	BA	977	G
4	BA	983	U
4	BA	986	U
4	BA	987	G
4	BA	992	U
4	BA	1015	C
4	BA	1016	C
4	BA	1018	U
4	BA	1026	A
4	BA	1027	G
4	BA	1028	A
4	BA	1035	A
4	BA	1038	A
4	BA	1041	A
4	BA	1045	U
4	BA	1050	C
4	BA	1055	A
4	BA	1056	G
4	BA	1057	A
4	BA	1063	U
4	BA	1064	G
4	BA	1071	U
4	BA	1072	A
4	BA	1074	U
4	BA	1075	U
4	BA	1081	U
4	BA	1089	G
4	BA	1091	U
4	BA	1092	A
4	BA	1093	A
4	BA	1127	A
4	BA	1129	C
4	BA	1138	G
4	BA	1140	A
4	BA	1146	G

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Mol	Chain	Res	Type
4	BA	1167	C
4	BA	1168	A
4	BA	1169	A
4	BA	1180	G
4	BA	1183	U
4	BA	1188	U
4	BA	1194	U
4	BA	1195	A
4	BA	1204	A
4	BA	1206	U
4	BA	1207	C
4	BA	1217	A
4	BA	1218	U
4	BA	1219	A
4	BA	1220	A
4	BA	1221	C
4	BA	1222	A
4	BA	1226	C
4	BA	1231	U
4	BA	1233	A
4	BA	1237	A
4	BA	1238	A
4	BA	1239	A
4	BA	1240	A
4	BA	1241	U
4	BA	1242	U
4	BA	1246	A
4	BA	1247	U
4	BA	1248	C
4	BA	1249	A
4	BA	1252	G
4	BA	1253	G
4	BA	1254	A
4	BA	1255	A
4	BA	1258	A
4	BA	1264	C
4	BA	1270	G
4	BA	1271	A
4	BA	1287	G
4	BA	1288	U
4	BA	1291	U
4	BA	1292	A

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Mol	Chain	Res	Type
4	BA	1294	A
4	BA	1298	C
4	BA	1299	C
4	BA	1314	U
4	BA	1321	C
4	BA	1323	U
4	BA	1325	G
4	BA	1326	A
4	BA	1328	G
4	BA	1332	G
4	BA	1336	A
4	BA	1341	A
4	BA	1342	C
4	BA	1352	G
4	BA	1358	G
4	BA	1385	U
4	BA	1386	U
4	BA	1387	A
4	BA	1388	A
4	BA	1389	A
4	BA	1390	G
4	BA	1395	A
4	BA	1396	C
4	BA	1405	A
4	BA	1409	C
4	BA	1420	A
4	BA	1425	A
4	BA	1426	G
4	BA	1429	C
4	BA	1433	U
4	BA	1436	U
4	BA	1438	U
4	BA	1445	U
4	BA	1448	A
4	BA	1457	A
4	BA	1468	A
4	BA	1474	G
4	BA	1476	A
4	BA	1493	A
4	BA	1494	A
4	BA	1498	C
4	BA	1504	A

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Mol	Chain	Res	Type
4	BA	1505	G
4	BA	1509	U
4	BA	1514	A
4	BA	1518	A
4	BA	1521	U
4	BA	1522	A
4	BA	1526	U
4	BA	1527	U
4	BA	1528	A
4	BA	1529	A
4	BA	1531	C
4	BA	1532	U
4	BA	1533	A
4	BA	1536	U
4	BA	1540	U
4	BA	1548	A
4	BA	1549	A
4	BA	1551	C
4	BA	1552	C
4	BA	1553	A
4	BA	1554	G
4	BA	1558	U
4	BA	1559	A
4	BA	1571	A
5	BB	3	U
5	BB	5	A
5	BB	7	G
5	BB	8	U
5	BB	9	A
5	BB	13	U
5	BB	21	C
5	BB	34	U
5	BB	43	C
5	BB	44	U
5	BB	46	G
5	BB	47	A
5	BB	48	U
5	BB	69	C
5	BB	70	A
47	AA	5	A
47	AA	10	U
47	AA	27	U

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Mol	Chain	Res	Type
47	AA	34	U
47	AA	42	A
47	AA	43	U
47	AA	44	U
47	AA	54	A
47	AA	58	U
47	AA	65	C
47	AA	67	C
47	AA	75	A
47	AA	81	U
47	AA	83	C
47	AA	102	G
47	AA	104	A
47	AA	115	A
47	AA	120	A
47	AA	139	A
47	AA	147	G
47	AA	152	A
47	AA	161	C
47	AA	164	C
47	AA	165	G
47	AA	170	A
47	AA	171	C
47	AA	173	G
47	AA	175	A
47	AA	186	U
47	AA	191	C
47	AA	194	U
47	AA	212	A
47	AA	216	A
47	AA	217	U
47	AA	219	U
47	AA	222	A
47	AA	223	U
47	AA	224	U
47	AA	231	G
47	AA	235	A
47	AA	238	C
47	AA	244	C
47	AA	247	G
47	AA	253	G
47	AA	256	G

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Mol	Chain	Res	Type
47	AA	257	U
47	AA	258	C
47	AA	273	A
47	AA	281	G
47	AA	287	C
47	AA	288	G
47	AA	294	A
47	AA	296	G
47	AA	297	A
47	AA	308	A
47	AA	309	A
47	AA	310	A
47	AA	314	A
47	AA	315	U
47	AA	316	C
47	AA	317	A
47	AA	320	A
47	AA	328	A
47	AA	337	C
47	AA	345	U
47	AA	354	C
47	AA	355	U
47	AA	361	A
47	AA	362	A
47	AA	368	A
47	AA	372	C
47	AA	395	C
47	AA	399	A
47	AA	407	U
47	AA	417	C
47	AA	421	A
47	AA	433	U
47	AA	455	A
47	AA	456	A
47	AA	457	C
47	AA	458	C
47	AA	461	A
47	AA	464	A
47	AA	471	U
47	AA	472	A
47	AA	477	A
47	AA	479	C

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Mol	Chain	Res	Type
47	AA	498	U
47	AA	502	A
47	AA	503	A
47	AA	518	A
47	AA	531	U
47	AA	532	G
47	AA	534	U
47	AA	538	C
47	AA	540	U
47	AA	541	C
47	AA	561	U
47	AA	566	U
47	AA	571	A
47	AA	574	C
47	AA	576	C
47	AA	579	A
47	AA	589	C
47	AA	593	C
47	AA	596	U
47	AA	597	U
47	AA	604	U
47	AA	616	C
47	AA	618	G
47	AA	619	C
47	AA	625	U
47	AA	628	G
47	AA	633	C
47	AA	639	A
47	AA	644	A
47	AA	645	A
47	AA	646	C
47	AA	648	A
47	AA	649	U
47	AA	652	U
47	AA	665	A
47	AA	681	A
47	AA	682	G
47	AA	696	U
47	AA	698	A
47	AA	707	A
47	AA	708	A
47	AA	709	A

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Mol	Chain	Res	Type
47	AA	711	A
47	AA	720	A
47	AA	722	A
47	AA	731	A
47	AA	732	U
47	AA	739	A
47	AA	743	A
47	AA	746	A
47	AA	766	C
47	AA	775	A
47	AA	790	A
47	AA	791	G
47	AA	802	A
47	AA	803	U
47	AA	807	G
47	AA	808	U
47	AA	823	G
47	AA	825	C
47	AA	838	A
47	AA	841	C
47	AA	842	A
47	AA	869	A
47	AA	870	G
47	AA	876	A
47	AA	877	A
47	AA	883	C
47	AA	885	U
47	AA	886	A
47	AA	893	A
47	AA	894	U
47	AA	897	C
47	AA	899	C
47	AA	902	C
47	AA	910	G
47	AA	917	C
47	AA	918	A
47	AA	919	A
47	AA	920	G
47	AA	923	G
47	AA	925	A
47	AA	928	A
47	AA	930	G

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Mol	Chain	Res	Type
47	AA	932	U
47	AA	943	G
47	AA	946	A
47	AA	955	G
47	AA	956	G
47	AA	959	U
47	AA	960	A
64	AV	2	G
64	AV	6	G
64	AV	8	U
64	AV	9	C
64	AV	14	A
64	AV	15	A
64	AV	16	A
64	AV	17	U
64	AV	18	A
64	AV	43	A
64	AV	44	U
64	AV	45	G
64	AV	51	U
64	AV	54	A
64	AV	55	C
64	AV	56	C
64	AV	58	U
64	AV	61	C
64	AV	63	G
64	AV	64	U
64	AV	65	A
64	AV	71	A
65	AX	3	A

All (3) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
4	BA	48	U
4	BA	1241	U
65	AX	2	A

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 347 ligands modelled in this entry, 341 are monoatomic - leaving 6 for Mogul analysis.

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	$\# Z > 2$
92	SPM	BA	3213	-	13,13,13	0.33	0	12,12,12	0.79	0
91	5GP	BA	3211	-	26,26,26	0.99	2 (7%)	40,40,40	1.73	8 (20%)
92	SPM	AA	3001	-	13,13,13	0.34	0	12,12,12	0.79	0
93	GTP	Ag	500	90	30,34,34	0.83	1 (3%)	46,54,54	1.77	11 (23%)
91	5GP	BA	3212	90	26,26,26	0.98	2 (7%)	40,40,40	1.68	8 (20%)
92	SPM	BA	3214	-	13,13,13	0.35	0	12,12,12	0.77	0

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
92	SPM	BA	3213	-	-	3/11/11/11	-
91	5GP	BA	3211	-	-	1/10/26/26	0/3/3/3
92	SPM	AA	3001	-	-	2/11/11/11	-
93	GTP	Ag	500	90	-	2/22/38/38	0/3/3/3
91	5GP	BA	3212	90	-	1/10/26/26	0/3/3/3
92	SPM	BA	3214	-	-	3/11/11/11	-

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
91	BA	3211	5GP	C6-N1	-2.55	1.34	1.38
91	BA	3212	5GP	C6-N1	-2.50	1.34	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
93	Ag	500	GTP	C2-N3	2.25	1.38	1.33
91	BA	3211	5GP	C5-N7	-2.06	1.35	1.39
91	BA	3212	5GP	C5-N7	-2.01	1.35	1.39

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
91	BA	3211	5GP	C5-C4-N3	-5.88	118.92	128.46
91	BA	3212	5GP	C5-C4-N3	-5.59	119.40	128.46
93	Ag	500	GTP	C5-C4-N3	-5.25	119.94	128.46
93	Ag	500	GTP	C2-N3-C4	4.61	120.52	112.30
91	BA	3211	5GP	C2-N3-C4	4.51	120.33	112.30
91	BA	3212	5GP	C2-N3-C4	4.35	120.05	112.30
93	Ag	500	GTP	PA-O3A-PB	-3.98	119.16	132.83
91	BA	3211	5GP	N9-C4-N3	3.75	133.46	125.94
93	Ag	500	GTP	N9-C4-N3	3.46	132.90	125.94
91	BA	3212	5GP	N9-C4-N3	3.42	132.80	125.94
93	Ag	500	GTP	PB-O3B-PG	-3.10	122.20	132.83
93	Ag	500	GTP	C2-N1-C6	-2.96	119.71	125.10
93	Ag	500	GTP	C5-C6-N1	2.56	119.69	113.19
91	BA	3211	5GP	C5-C6-N1	2.48	119.50	113.19
91	BA	3212	5GP	C5-C6-N1	2.47	119.47	113.19
93	Ag	500	GTP	C8-N7-C5	2.46	108.69	104.24
93	Ag	500	GTP	N9-C8-N7	-2.46	108.76	113.39
93	Ag	500	GTP	O6-C6-C5	-2.44	120.13	126.60
91	BA	3211	5GP	O6-C6-C5	-2.37	120.31	126.60
91	BA	3212	5GP	O6-C6-C5	-2.36	120.33	126.60
91	BA	3211	5GP	C2-N1-C6	-2.35	120.81	125.10
91	BA	3212	5GP	C2-N1-C6	-2.32	120.88	125.10
93	Ag	500	GTP	C3'-C2'-C1'	2.31	105.81	101.43
91	BA	3212	5GP	N9-C8-N7	-2.27	109.12	113.39
91	BA	3211	5GP	N9-C8-N7	-2.22	109.21	113.39
91	BA	3212	5GP	C4-C5-N7	-2.12	107.36	110.72
91	BA	3211	5GP	C4-C5-N7	-2.05	107.47	110.72

There are no chirality outliers.

All (12) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
92	BA	3213	SPM	C7-C8-C9-N10
92	BA	3214	SPM	N5-C6-C7-C8
93	Ag	500	GTP	O4'-C4'-C5'-O5'

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Mol	Chain	Res	Type	Atoms
93	Ag	500	GTP	C3'-C4'-C5'-O5'
92	AA	3001	SPM	C6-C7-C8-C9
92	BA	3214	SPM	C7-C8-C9-N10
92	BA	3214	SPM	C8-C9-N10-C11
92	AA	3001	SPM	C8-C9-N10-C11
91	BA	3212	5GP	C5'-O5'-P-OP2
92	BA	3213	SPM	C7-C6-N5-C4
91	BA	3211	5GP	O4'-C4'-C5'-O5'
92	BA	3213	SPM	C6-C7-C8-C9

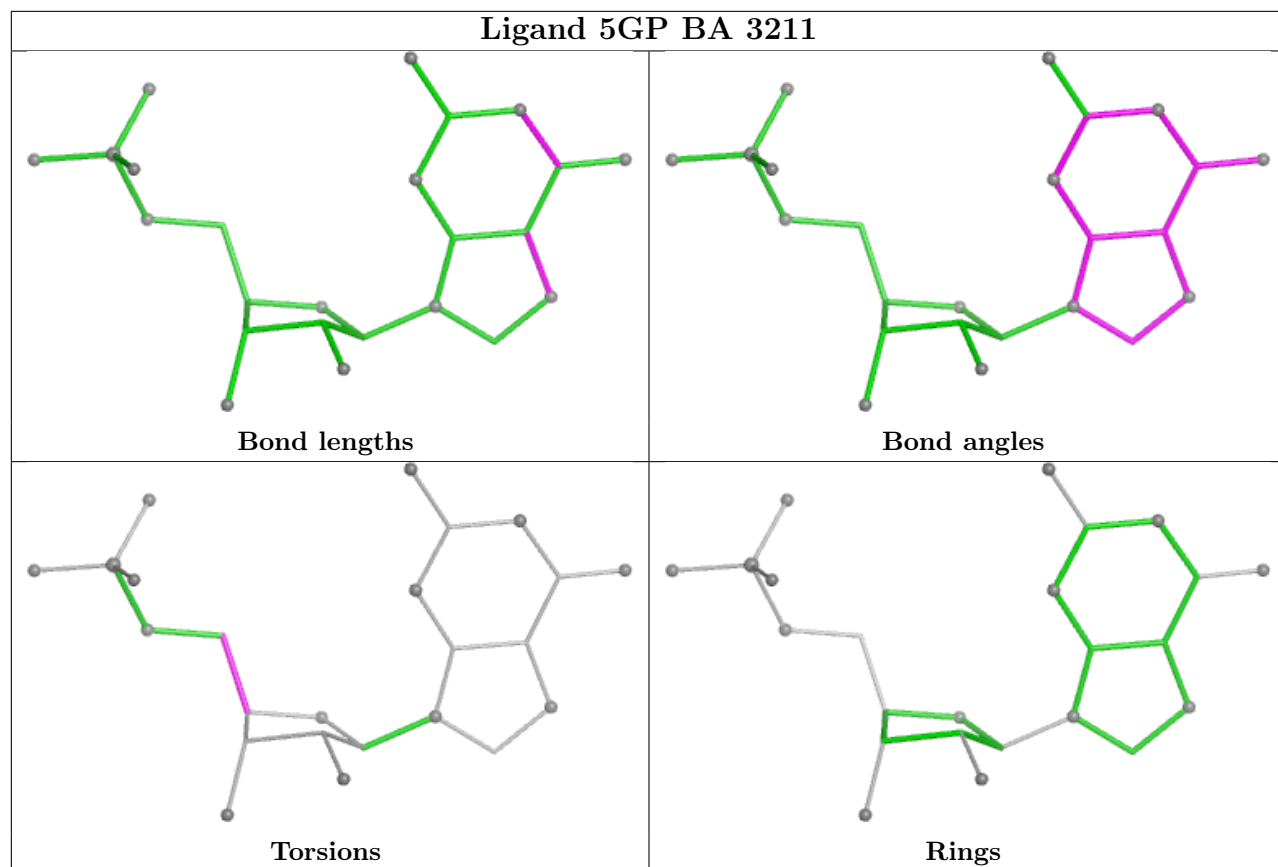
There are no ring outliers.

4 monomers are involved in 8 short contacts:

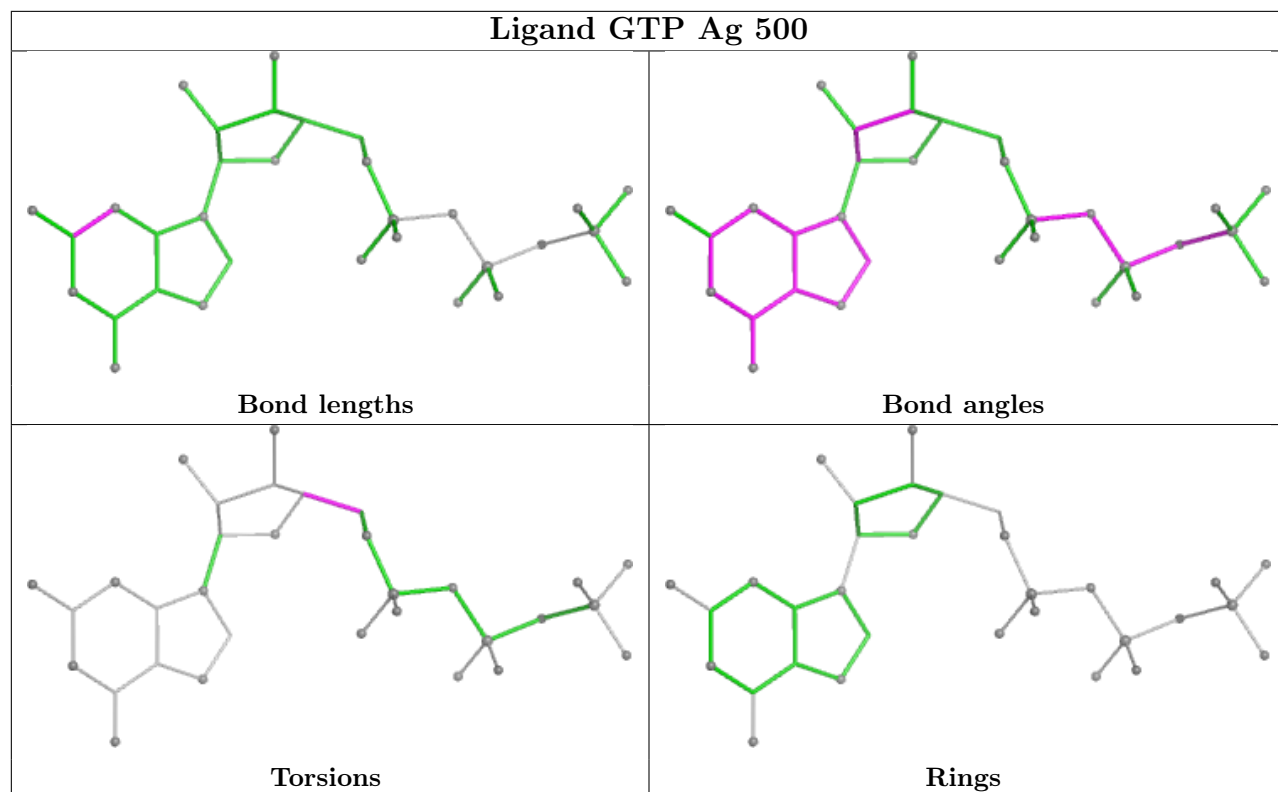
Mol	Chain	Res	Type	Clashes	Symm-Clashes
92	BA	3213	SPM	2	0
92	AA	3001	SPM	1	0
93	Ag	500	GTP	3	0
91	BA	3212	5GP	2	0

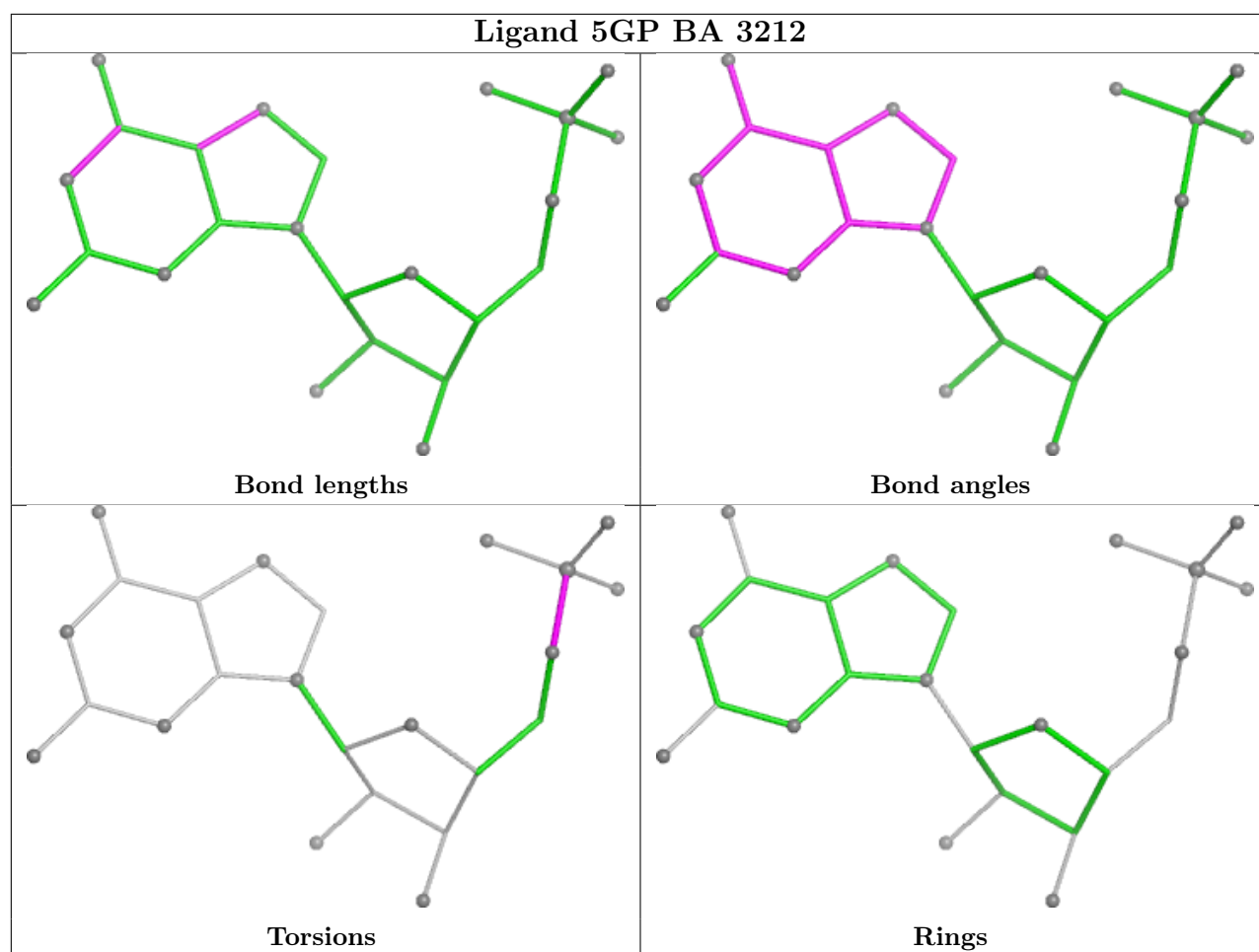
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

Ligand 5GP BA 3211



Ligand GTP Ag 500





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

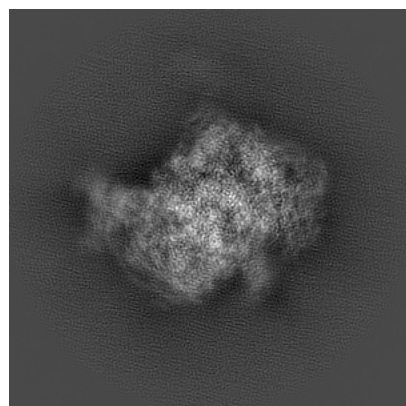
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-12529. These allow visual inspection of the internal detail of the map and identification of artifacts.

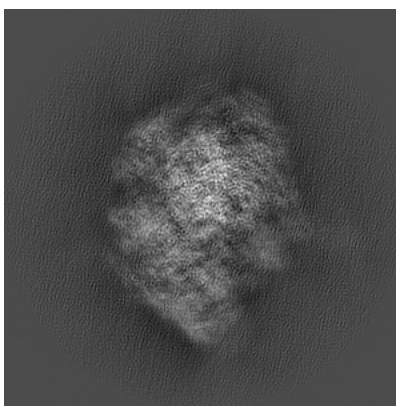
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

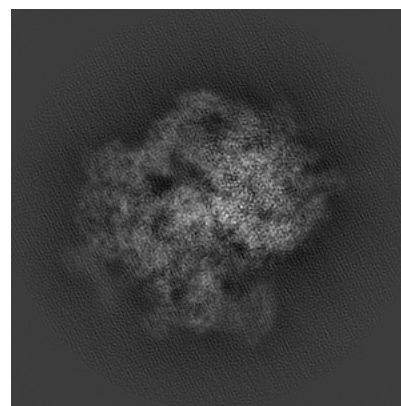
6.1.1 Primary map



X

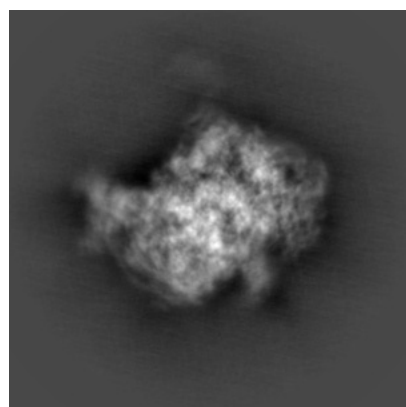


Y

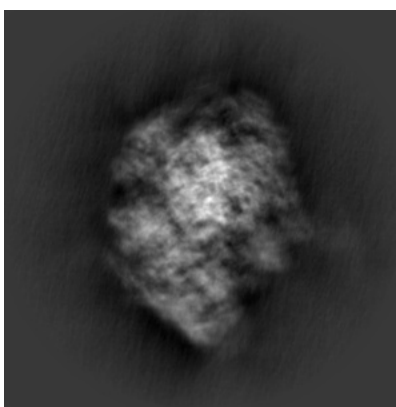


Z

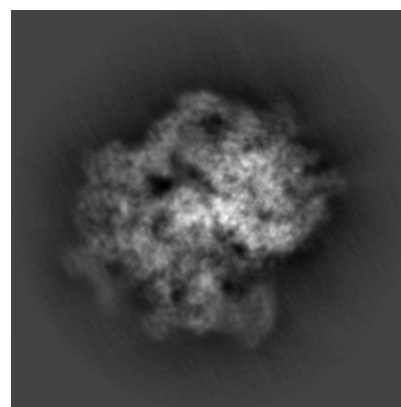
6.1.2 Raw map



X



Y

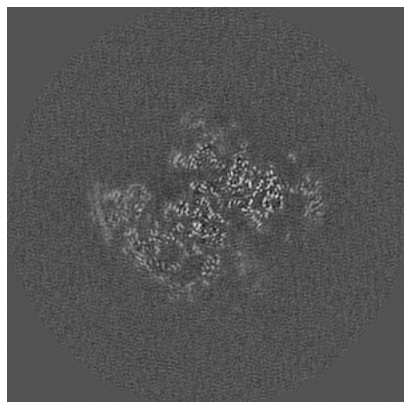


Z

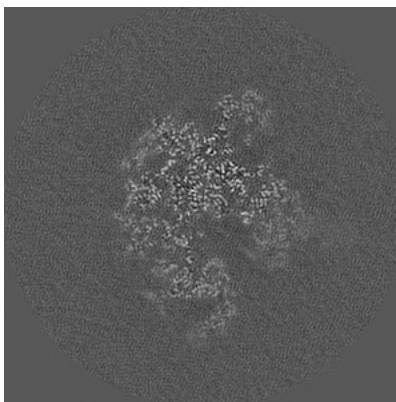
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

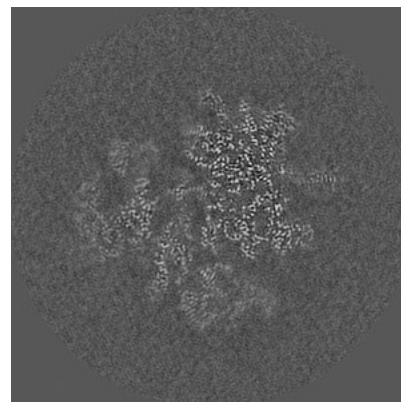
6.2.1 Primary map



X Index: 160

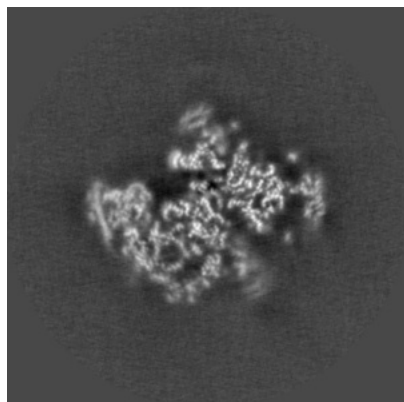


Y Index: 160

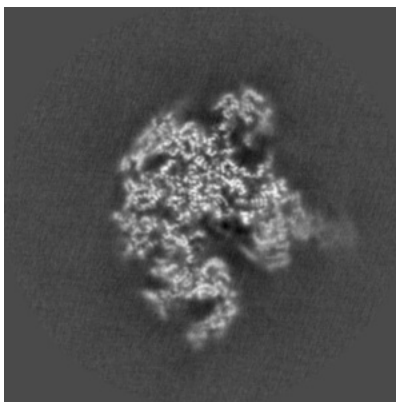


Z Index: 160

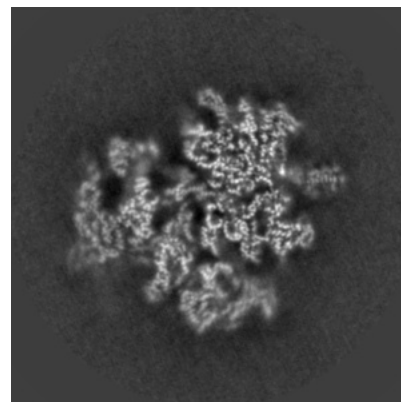
6.2.2 Raw map



X Index: 160



Y Index: 160

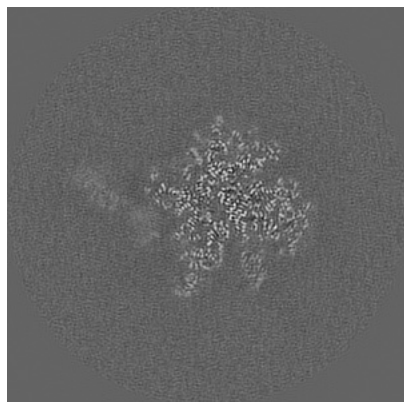


Z Index: 160

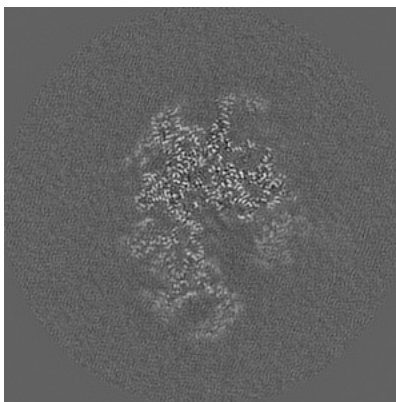
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

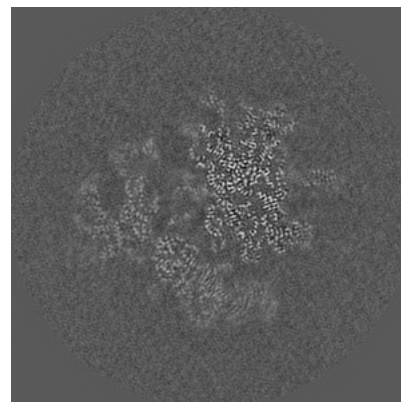
6.3.1 Primary map



X Index: 184

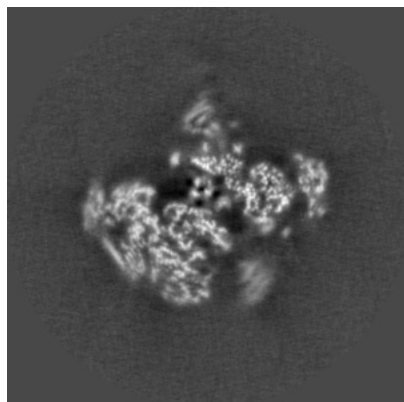


Y Index: 164

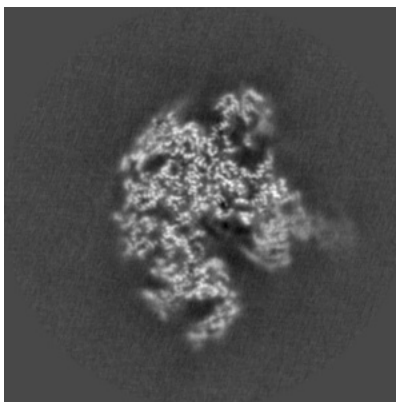


Z Index: 163

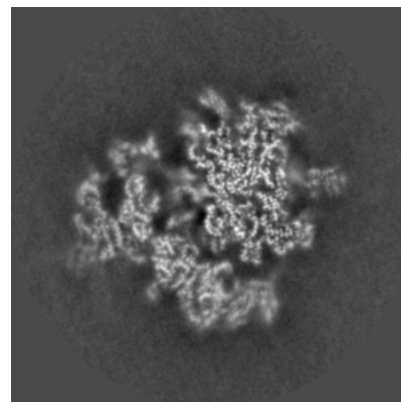
6.3.2 Raw map



X Index: 154



Y Index: 161

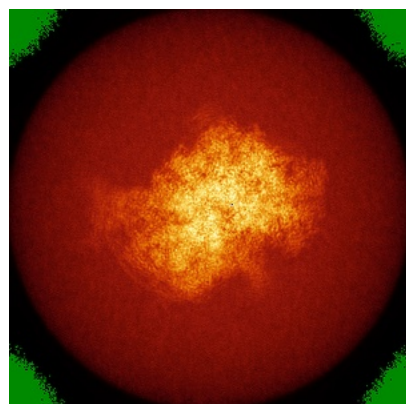


Z Index: 163

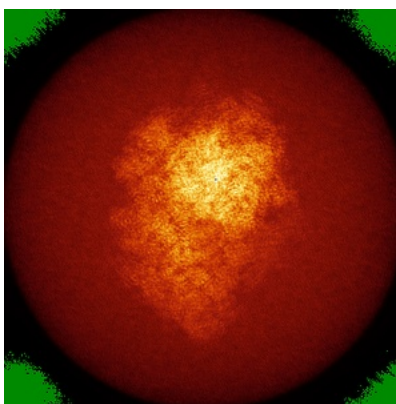
The images above show the largest variance slices of the map in three orthogonal directions.

6.4 Orthogonal standard-deviation projections (False-color) [i](#)

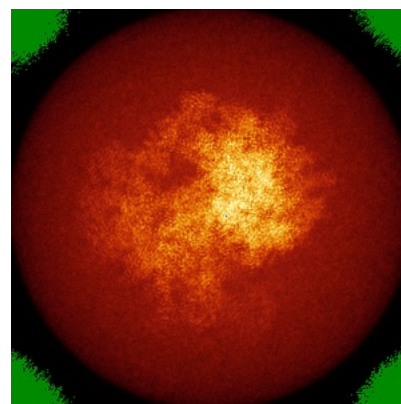
6.4.1 Primary map



X

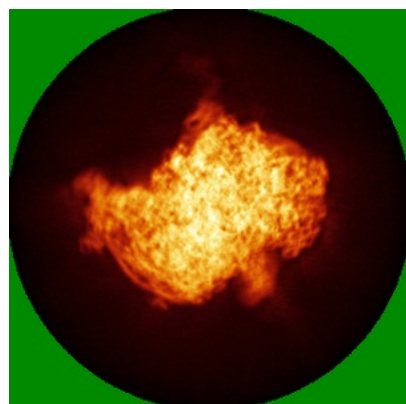


Y

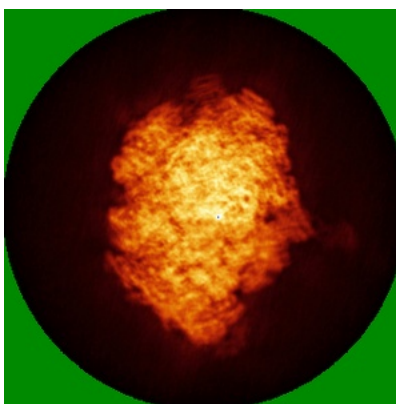


Z

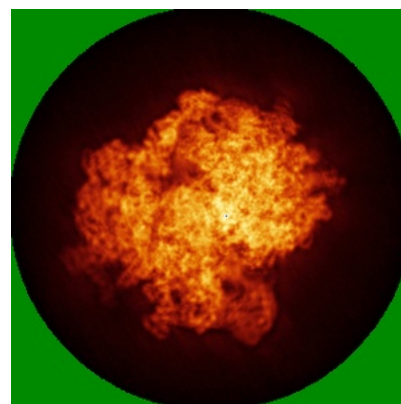
6.4.2 Raw map



X



Y

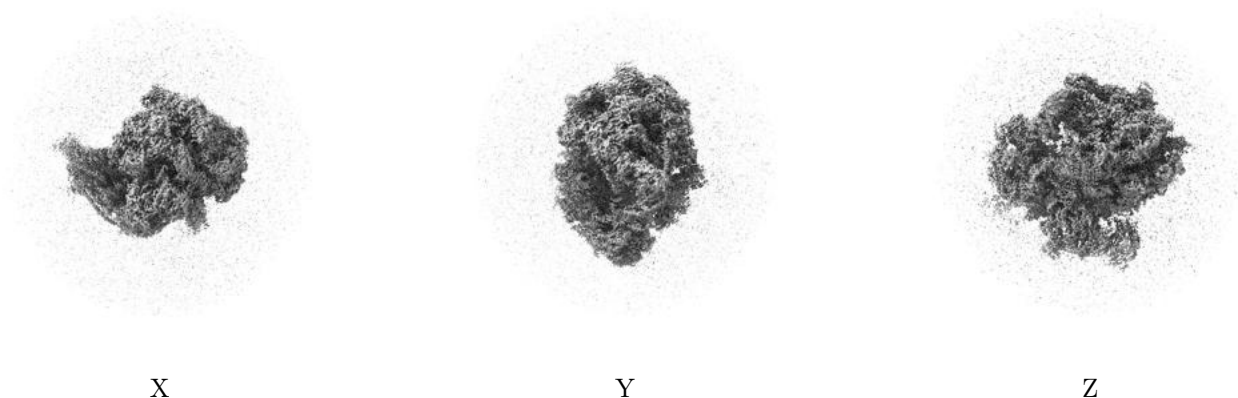


Z

The images above show the map standard deviation projections with false color in three orthogonal directions. Minimum values are shown in green, max in blue, and dark to light orange shades represent small to large values respectively.

6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.13. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.5.2 Raw map



These images show the 3D surface of the raw map. The raw map's contour level was selected so that its surface encloses the same volume as the primary map does at its recommended contour level.

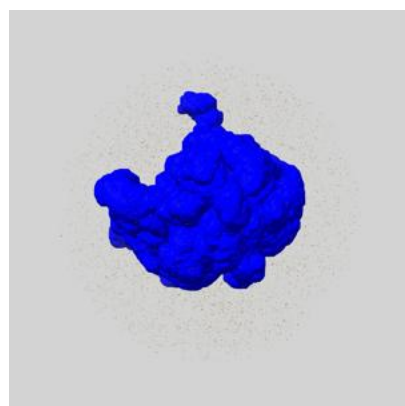
6.6 Mask visualisation [i](#)

This section shows the 3D surface view of the primary map at 50% transparency overlaid with the specified mask at 0% transparency

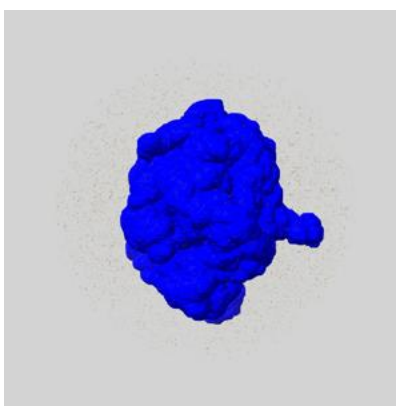
A mask typically either:

- Encompasses the whole structure
- Separates out a domain, a functional unit, a monomer or an area of interest from a larger structure

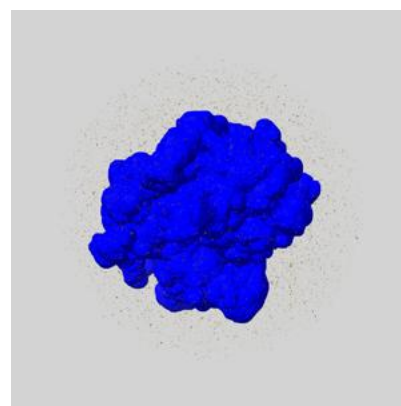
6.6.1 emd_12529_msk_1.map [i](#)



X



Y

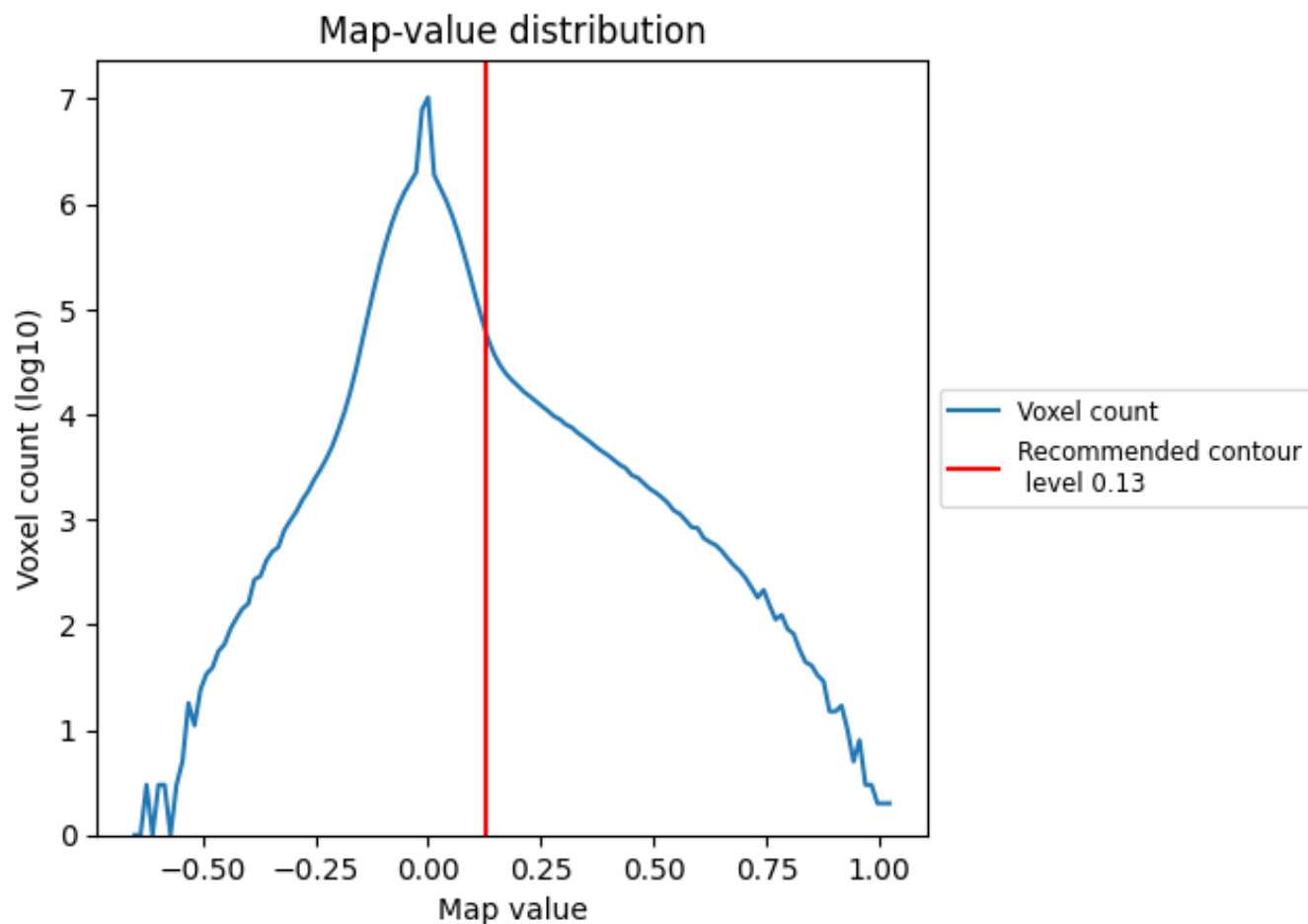


Z

7 Map analysis [i](#)

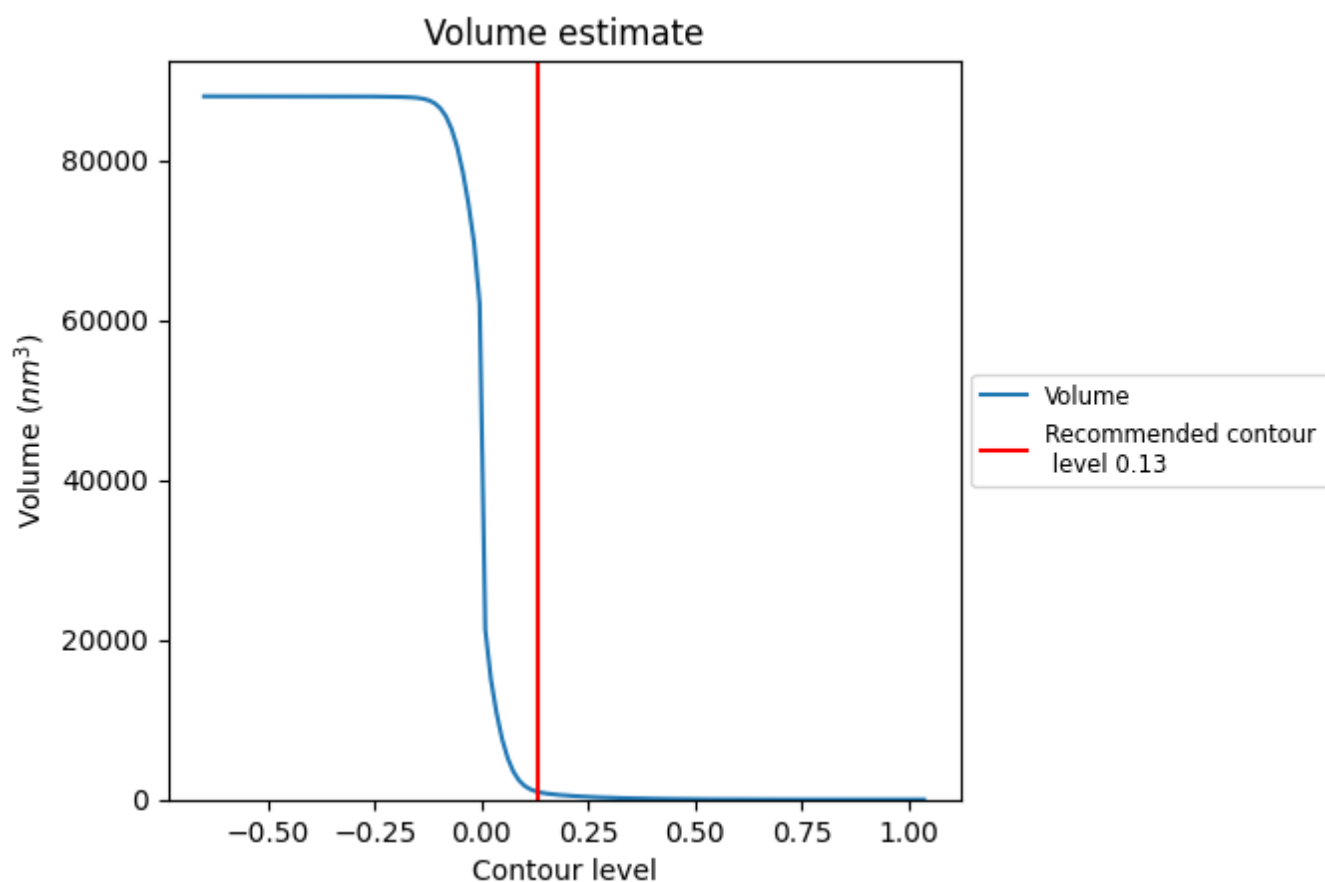
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



The map-value distribution is plotted in 128 intervals along the x-axis. The y-axis is logarithmic. A spike in this graph at zero usually indicates that the volume has been masked.

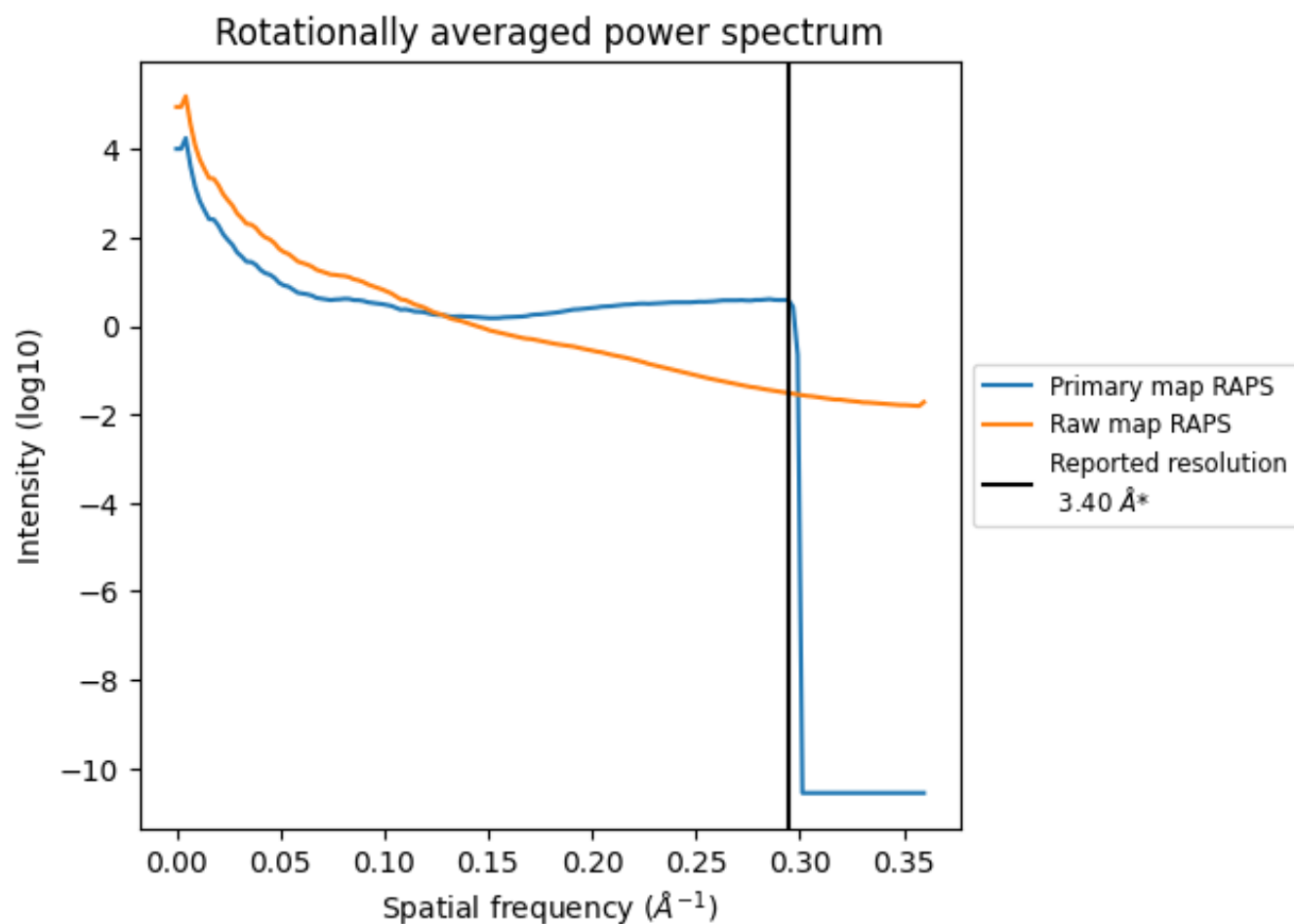
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 992 nm³; this corresponds to an approximate mass of 897 kDa.

The volume estimate graph shows how the enclosed volume varies with the contour level. The recommended contour level is shown as a vertical line and the intersection between the line and the curve gives the volume of the enclosed surface at the given level.

7.3 Rotationally averaged power spectrum ⓘ

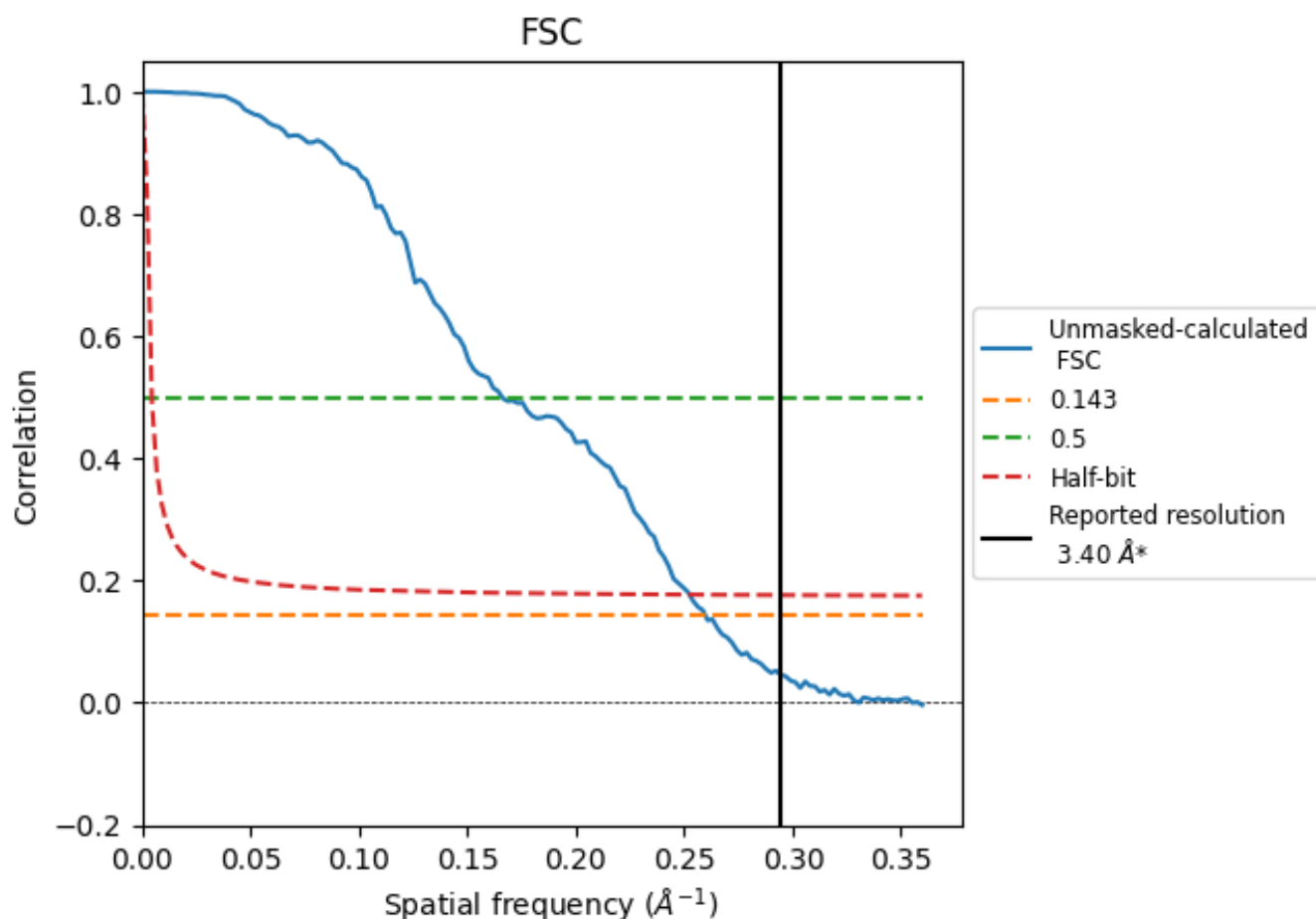


*Reported resolution corresponds to spatial frequency of 0.294 Å⁻¹

8 Fourier-Shell correlation [i](#)

Fourier-Shell Correlation (FSC) is the most commonly used method to estimate the resolution of single-particle and subtomogram-averaged maps. The shape of the curve depends on the imposed symmetry, mask and whether or not the two 3D reconstructions used were processed from a common reference. The reported resolution is shown as a black line. A curve is displayed for the half-bit criterion in addition to lines showing the 0.143 gold standard cut-off and 0.5 cut-off.

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of 0.294 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

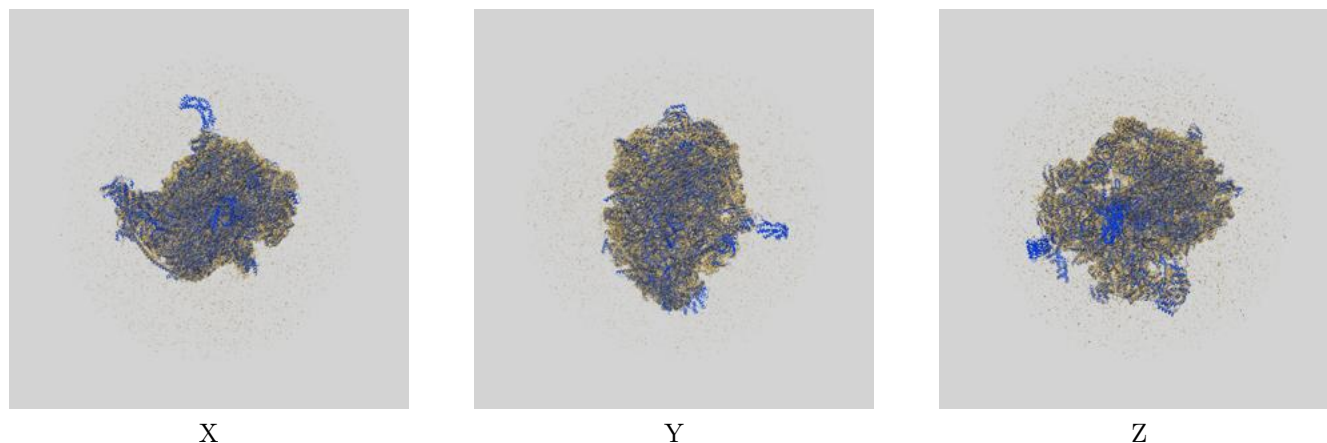
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.40	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	3.85	6.02	3.96

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 3.85 differs from the reported value 3.4 by more than 10 %

9 Map-model fit [i](#)

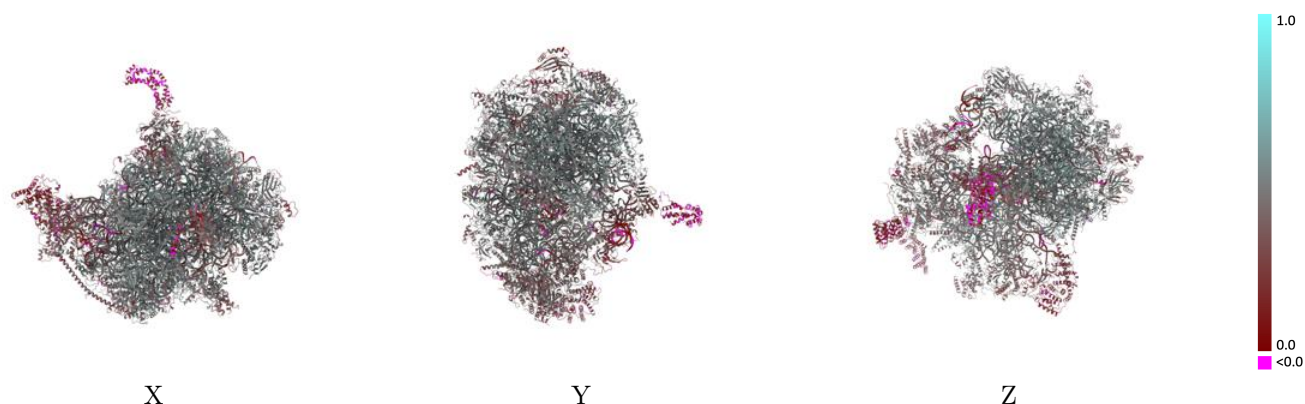
This section contains information regarding the fit between EMDB map EMD-12529 and PDB model 7NQL. Per-residue inclusion information can be found in section 3 on page 25.

9.1 Map-model overlay [i](#)



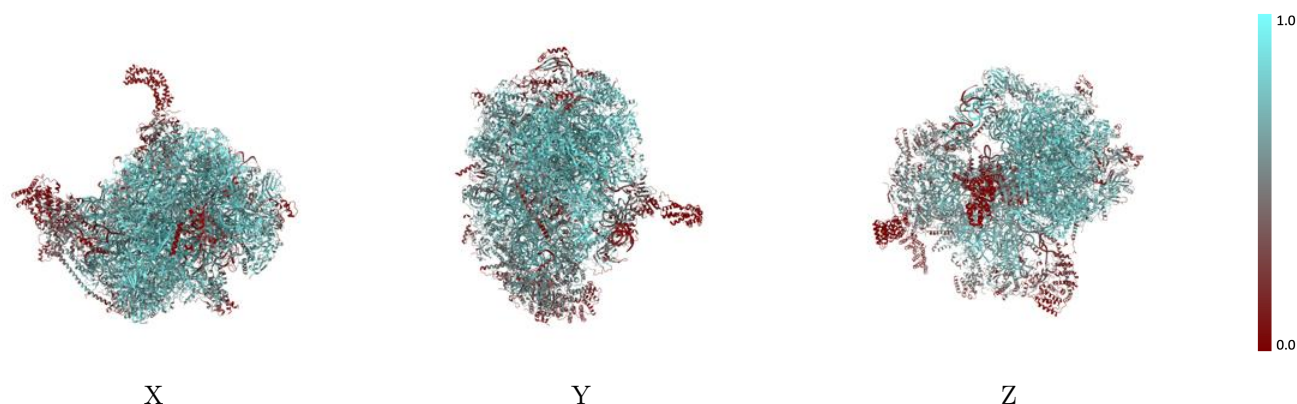
The images above show the 3D surface view of the map at the recommended contour level 0.13 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

9.2 Q-score mapped to coordinate model [i](#)



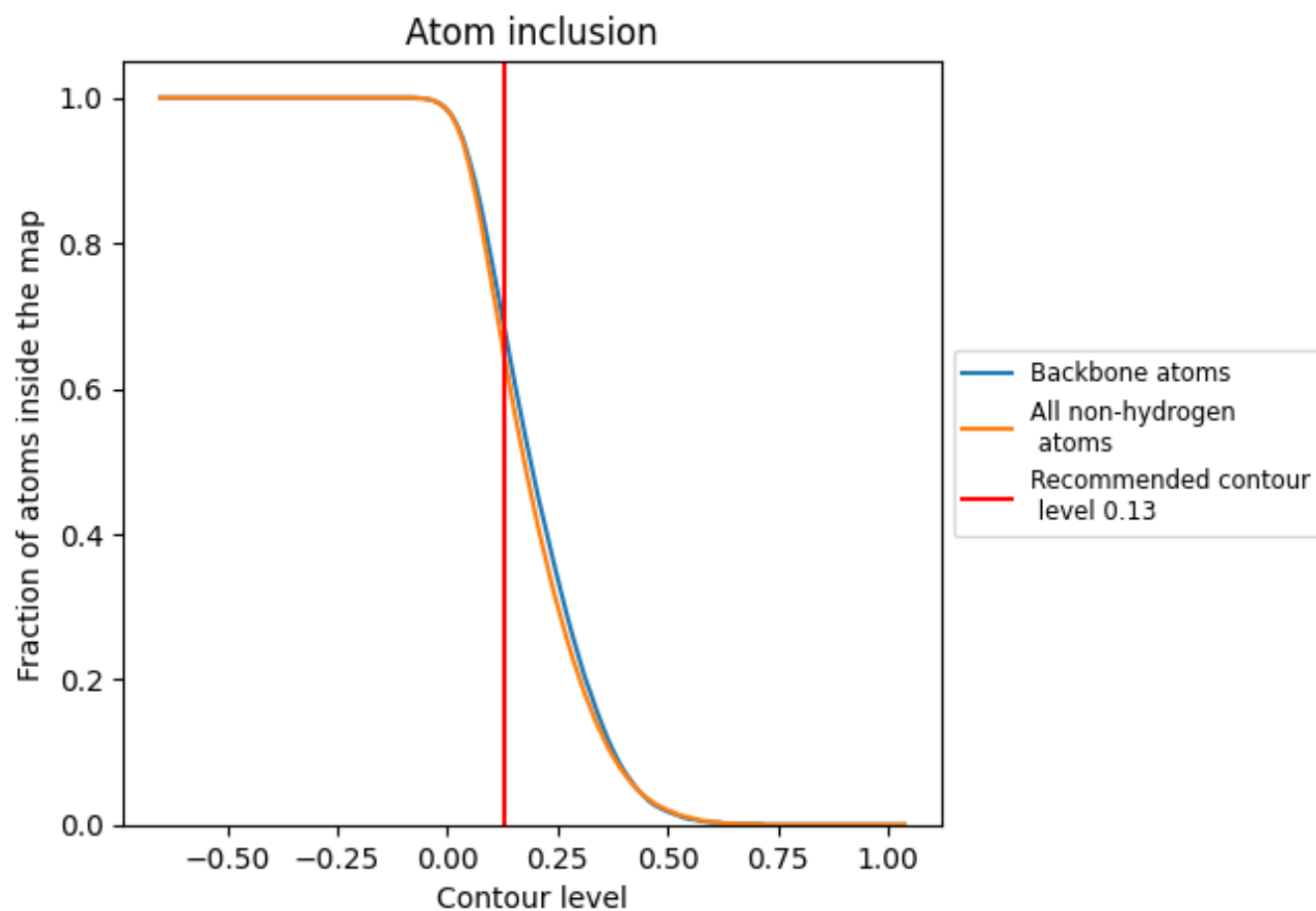
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions at which atoms can be resolved. Low Q-score values may therefore be expected for many entries.

9.3 Atom inclusion mapped to coordinate model [i](#)



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (0.13).

9.4 Atom inclusion [i](#)



At the recommended contour level, 68% of all backbone atoms, 64% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary ⓘ

The table lists the average atom inclusion at the recommended contour level (0.13) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.6420	 0.4370
AA	 0.7650	 0.4380
AB	 0.6840	 0.4670
AC	 0.6190	 0.4730
AE	 0.5570	 0.4510
AF	 0.6700	 0.4750
AG	 0.5490	 0.4250
AI	 0.5230	 0.4100
AJ	 0.5500	 0.4190
AK	 0.6700	 0.4770
AL	 0.6400	 0.4840
AN	 0.6800	 0.4640
AO	 0.5520	 0.4330
AP	 0.5010	 0.3880
AQ	 0.6410	 0.4620
AR	 0.6560	 0.4760
AU	 0.7080	 0.4840
AV	 0.6020	 0.3620
AX	 0.7580	 0.4820
AZ	 0.3330	 0.2130
Aa	 0.3800	 0.3680
Ab	 0.4600	 0.4010
Ac	 0.6090	 0.4400
Ad	 0.4730	 0.3740
Ae	 0.1410	 0.2260
Af	 0.5650	 0.4550
Ag	 0.4680	 0.3730
Ah	 0.4130	 0.3570
Ai	 0.4960	 0.4080
Aj	 0.3000	 0.3350
Ak	 0.4400	 0.3790
Am	 0.5490	 0.4370
An	 0.6750	 0.5050
Ao	 0.0780	 0.2060
Ap	 0.5070	 0.3980



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Chain	Atom inclusion	Q-score
B0	 0.8310	 0.5370
B1	 0.5990	 0.4580
B2	 0.7020	 0.4790
B3	 0.8130	 0.5250
B4	 0.6220	 0.4240
B5	 0.7420	 0.5020
B6	 0.4350	 0.4340
B7	 0.8540	 0.5460
B8	 0.8430	 0.5460
B9	 0.8720	 0.5440
BA	 0.8320	 0.4810
BB	 0.5350	 0.2780
BD	 0.7830	 0.5180
BE	 0.7900	 0.5160
BF	 0.7670	 0.5020
BI	 0.4470	 0.4260
BJ	 0.2770	 0.3110
BK	 0.1230	 0.2410
BL	 0.7030	 0.4940
BN	 0.8280	 0.5350
BO	 0.7450	 0.5080
BP	 0.7560	 0.4990
BQ	 0.7740	 0.5150
BR	 0.7770	 0.5200
BS	 0.7680	 0.5040
BT	 0.6900	 0.4910
BU	 0.8020	 0.5210
BV	 0.7890	 0.5190
BW	 0.7590	 0.5210
BX	 0.6480	 0.4650
BY	 0.3520	 0.3860
Ba	 0.7320	 0.4820
Bb	 0.7320	 0.4660
Bc	 0.6420	 0.4460
Bd	 0.5270	 0.3800
Be	 0.5600	 0.4430
Bf	 0.6880	 0.4540
Bg	 0.7870	 0.5140
Bh	 0.6970	 0.4720
Bi	 0.2860	 0.3600
Bj	 0.4720	 0.3200
Bk	 0.5550	 0.4230

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Chain	Atom inclusion	Q-score
Bl	 0.7750	 0.5020
Bm	 0.2760	 0.3810
Bn	 0.7940	 0.5190
Bo	 0.7340	 0.4900
Bp	 0.3290	 0.3560
Bq	 0.3680	 0.3810
Bt	 0.7990	 0.5180
Bu	 0.4600	 0.3940
Bv	 0.4510	 0.4080
Bw	 0.7320	 0.4860
Bx	 0.7240	 0.4710
CL	 0.0090	 0.1130
DL	 0.0050	 0.1040
EL	 0.0000	 0.0740
FL	 0.0000	 0.0320
GL	 0.0000	 0.0010
HL	 0.0000	 0.0060