Full wwPDB EM Validation Report ⓘ

Jun 26, 2025 – 04:09 AM JST

PDB ID	:	8H9V / pdb_00008h9v
EMDB ID	:	EMD-34583
Title	:	Human ATP synthase state 3b (combined)
Authors	:	Lai, Y.; Zhang, Y.; Liu, F.; Gao, Y.; Gong, H.; Rao, Z.
Deposited on	:	2022-10-25
Resolution	:	3.02 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:

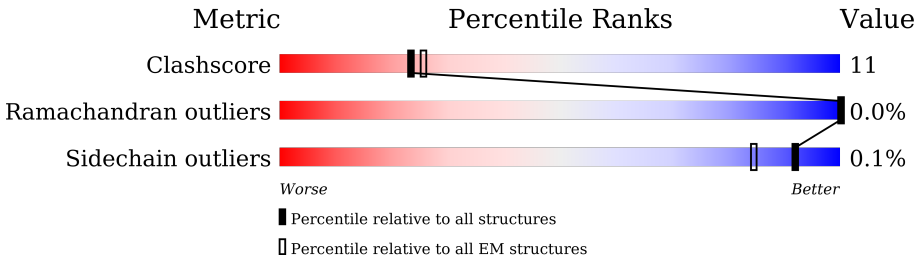
EMDB validation analysis	:	0.0.1.dev118
Mogul	:	1.8.5 (274361), CSD as541be (2020)
MolProbity	:	4-5-2 with Phenix2.0rc1
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
MapQ	:	1.9.13
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.44

1 Overall quality at a glance (i)

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

The reported resolution of this entry is 3.02 Å.

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



Metric	Whole archive (#Entries)	EM structures (#Entries)
Clashscore	210492	15764
Ramachandran outliers	207382	16835
Sidechain outliers	206894	16415

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 <=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	1	75	81% 17%
1	2	75	81% 19%
1	3	75	75% 25%
1	4	75	65% 35%
1	5	75	53% 47%
1	6	75	63% 37%
1	7	75	7% 79% 21%
1	8	75	80% 20%

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Mol	Chain	Length	Quality of chain
2	H	146	63% 27% 10%
3	I	51	71% 18% 12%
4	K	214	60% 33% 7%
5	M	160	59% 38%
6	N	226	69% 30%
7	P	58	53% 17% 29%
8	Q	68	49% 26% 25%
9	R	93	58% 22% 20%
10	S	102	44% 31% 25%
11	T	69	45% 19% 36%
12	L	108	38% 24% 38%
13	A	510	73% 23% 5%
13	B	510	76% 17% 6%
13	C	510	77% 15% 8%
14	D	482	6% 65% 24% 10%
14	E	482	74% 20% 5%
14	F	482	78% 18%
15	G	273	5% 78% 21%
16	O	190	58% 66% 33%

2 Entry composition

There are 19 unique types of molecules in this entry. The entry contains 37921 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 ATP synthase F(0) complex subunit C1, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	4	75	Total	C	N	O	S	0	0
			534	353	83	94	4		
1	5	75	Total	C	N	O	S	0	0
			534	353	83	94	4		
1	6	75	Total	C	N	O	S	0	0
			534	353	83	94	4		
1	7	75	Total	C	N	O	S	0	0
			534	353	83	94	4		
1	8	75	Total	C	N	O	S	0	0
			534	353	83	94	4		
1	1	75	Total	C	N	O	S	0	0
			534	353	83	94	4		
1	2	75	Total	C	N	O	S	0	0
			534	353	83	94	4		
1	3	75	Total	C	N	O	S	0	0
			534	353	83	94	4		

- Molecule 2 is a protein called ATP synthase subunit delta, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	H	132	Total	C	N	O	S	0	0
			975	614	164	195	2		

- Molecule 3 is a protein called ATP synthase subunit epsilon, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	I	45	Total	C	N	O	S	0	0
			354	225	64	64	1		

- Molecule 4 is a protein called ATP synthase F(0) complex subunit B1, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	K	198	Total	C	N	O	S	0	0
			1573	1013	270	285	5		

- Molecule 5 is a protein called ATP synthase subunit d, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	M	156	Total	C	N	O	S	0	0
			1259	813	199	243	4		

- Molecule 6 is a protein called ATP synthase subunit a.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	N	223	Total	C	N	O	S	0	0
			1718	1145	273	289	11		

- Molecule 7 is a protein called ATP synthase subunit ATP5MJ, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
7	P	41	Total	C	N	O	S	0	0
			344	234	53	54	3		

- Molecule 8 is a protein called ATP synthase protein 8.

Mol	Chain	Residues	Atoms					AltConf	Trace
8	Q	51	Total	C	N	O	S	0	0
			422	281	67	68	6		

- Molecule 9 is a protein called ATP synthase subunit f, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
9	R	74	Total	C	N	O	S	0	0
			621	411	104	103	3		

- Molecule 10 is a protein called ATP synthase subunit g, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
10	S	77	Total	C	N	O	S	0	0
			605	401	96	107	1		

- Molecule 11 is a protein called ATP synthase subunit e, mitochondrial.

Mol	Chain	Residues	Atoms				AltConf	Trace
11	T	44	Total	C	N	O	0	0
			346	224	63	59		

- Molecule 12 is a protein called ATP synthase-coupling factor 6, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
12	L	67	Total	C	N	O	S	0	0
			549	350	91	106	2		

- Molecule 13 is a protein called ATP synthase subunit alpha, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
13	A	486	Total	C	N	O	S	0	0
			3706	2334	655	705	12		
13	B	477	Total	C	N	O	S	0	0
			3639	2290	645	692	12		
13	C	470	Total	C	N	O	S	0	0
			3587	2261	636	678	12		

- Molecule 14 is a protein called ATP synthase subunit beta, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
14	E	456	Total	C	N	O	S	0	0
			3458	2192	588	666	12		
14	F	466	Total	C	N	O	S	0	0
			3529	2238	598	680	13		
14	D	432	Total	C	N	O	S	0	0
			3272	2074	561	625	12		

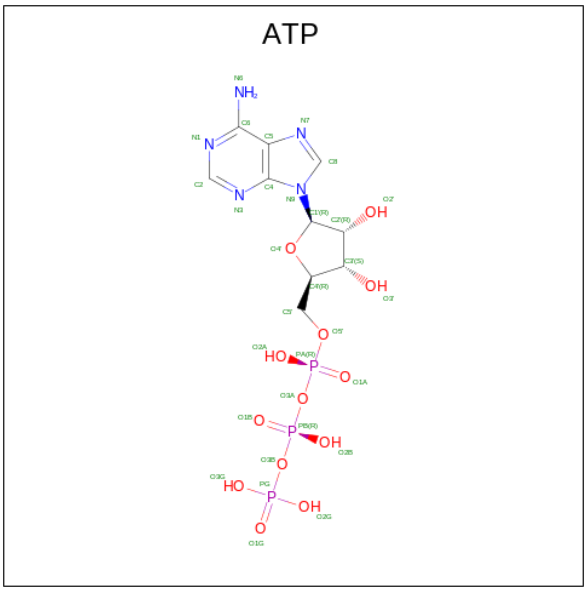
- Molecule 15 is a protein called ATP synthase subunit gamma, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
15	G	271	Total	C	N	O	S	0	0
			2103	1329	359	406	9		

- Molecule 16 is a protein called ATP synthase subunit O, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
16	O	187	Total	C	N	O	S	0	0
			1437	909	252	270	6		

- Molecule 17 is ADENOSINE-5'-TRIPHOSPHATE (CCD ID: ATP) (formula: $C_{10}H_{16}N_5O_{13}P_3$) (labeled as "Ligand of Interest" by depositor).

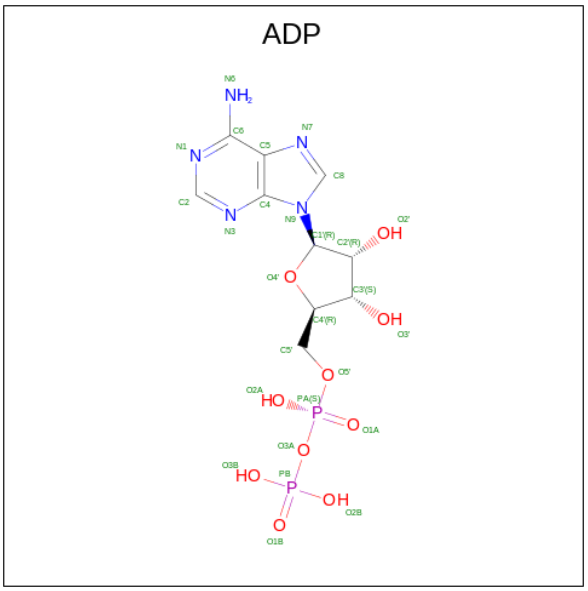


Mol	Chain	Residues	Atoms					AltConf
17	A	1	Total	C	N	O	P	0
			31	10	5	13	3	
17	B	1	Total	C	N	O	P	0
			31	10	5	13	3	
17	C	1	Total	C	N	O	P	0
			31	10	5	13	3	

- Molecule 18 is MAGNESIUM ION (CCD ID: MG) (formula: Mg) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		AltConf
18	A	1	Total	Mg	0
			1	1	
18	B	1	Total	Mg	0
			1	1	
18	C	1	Total	Mg	0
			1	1	
18	F	1	Total	Mg	0
			1	1	
18	D	1	Total	Mg	0
			1	1	

- Molecule 19 is ADENOSINE-5'-DIPHOSPHATE (CCD ID: ADP) (formula: $C_{10}H_{15}N_5O_{10}P_2$) (labeled as "Ligand of Interest" by depositor).



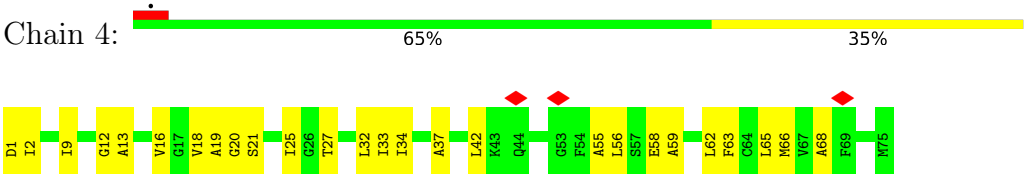
Mol	Chain	Residues	Atoms					AltConf
19	F	1	Total	C	N	O	P	0
			27	10	5	10	2	
19	D	1	Total	C	N	O	P	0
			27	10	5	10	2	

3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion in map density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red diamond above a residue indicates a poor fit to the EM map for this residue (all-atom inclusion < 40%). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

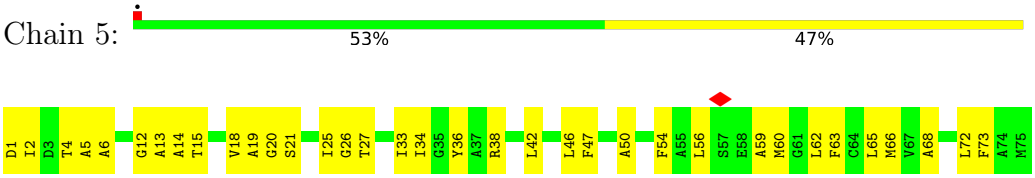
• Molecule 1: ATP synthase F(0) complex subunit C1, mitochondrial

Chain 4:



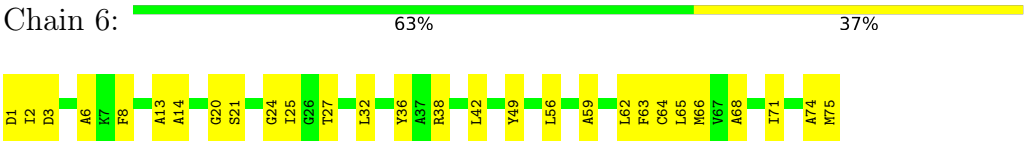
• Molecule 1: ATP synthase F(0) complex subunit C1, mitochondrial

Chain 5:



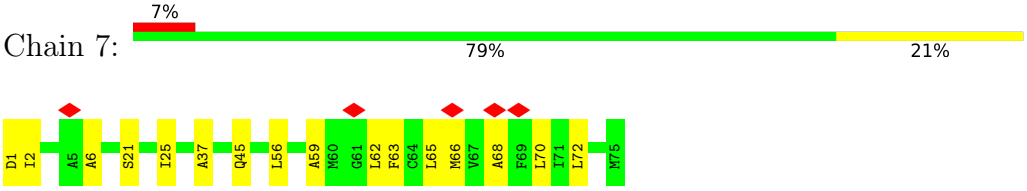
• Molecule 1: ATP synthase F(0) complex subunit C1, mitochondrial

Chain 6:



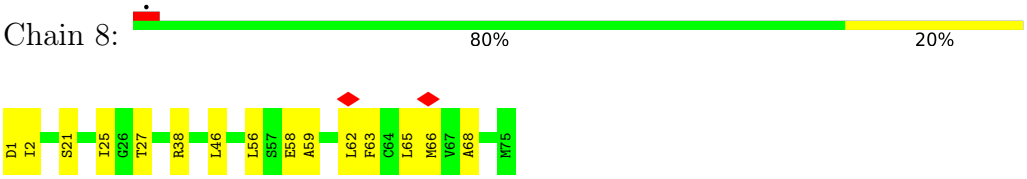
• Molecule 1: ATP synthase F(0) complex subunit C1, mitochondrial

Chain 7:



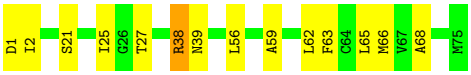
• Molecule 1: ATP synthase F(0) complex subunit C1, mitochondrial

Chain 8:



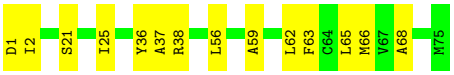
• Molecule 1: ATP synthase F(0) complex subunit C1, mitochondrial

Chain 1:



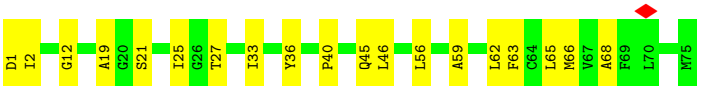
• Molecule 1: ATP synthase F(0) complex subunit C1, mitochondrial

Chain 2:



• Molecule 1: ATP synthase F(0) complex subunit C1, mitochondrial

Chain 3:

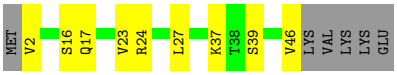


• Molecule 2: ATP synthase subunit delta, mitochondrial

Chain H:

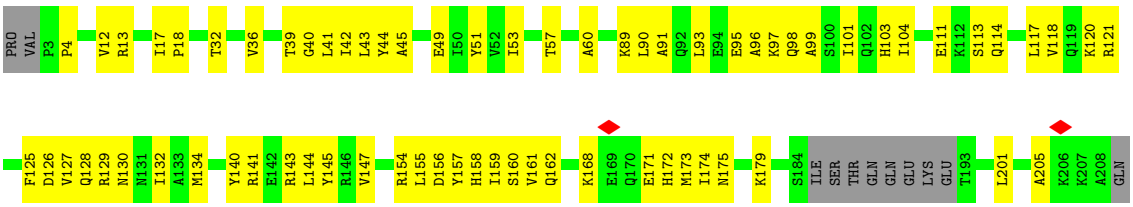
• Molecule 3: ATP synthase subunit epsilon, mitochondrial

Chain I:



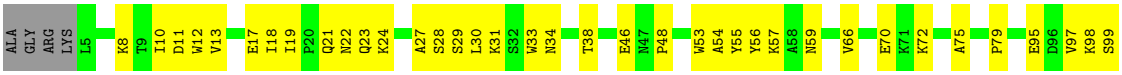
• Molecule 4: ATP synthase F(0) complex subunit B1, mitochondrial

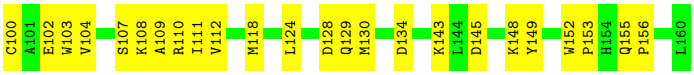
Chain K:



• Molecule 5: ATP synthase subunit d, mitochondrial

Chain M:



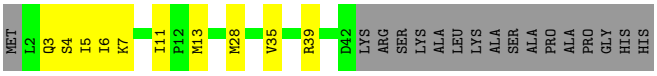


• Molecule 6: ATP synthase subunit a

Chain N:

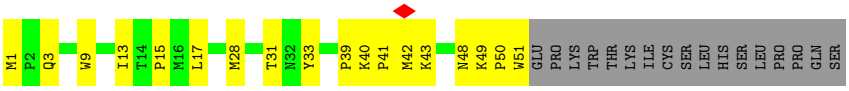
• Molecule 7: ATP synthase subunit ATP5MJ, mitochondrial

Chain P:



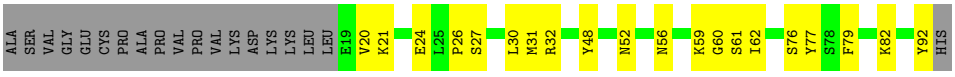
• Molecule 8: ATP synthase protein 8

Chain Q:



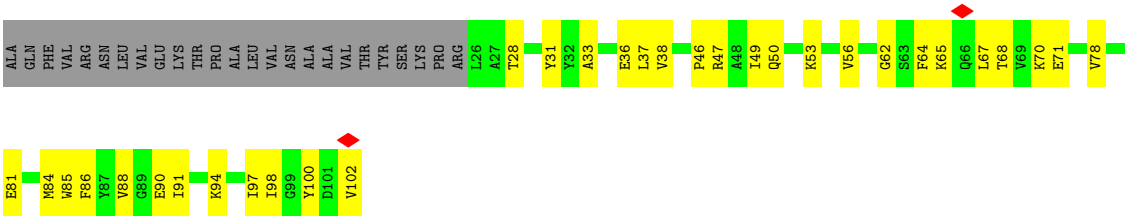
• Molecule 9: ATP synthase subunit f, mitochondrial

Chain R:



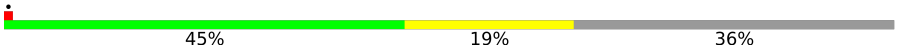
• Molecule 10: ATP synthase subunit g, mitochondrial

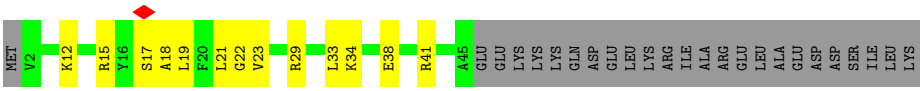
Chain S:



• Molecule 11: ATP synthase subunit e, mitochondrial

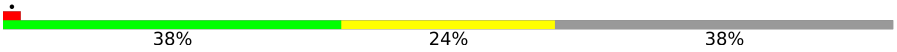
Chain T:

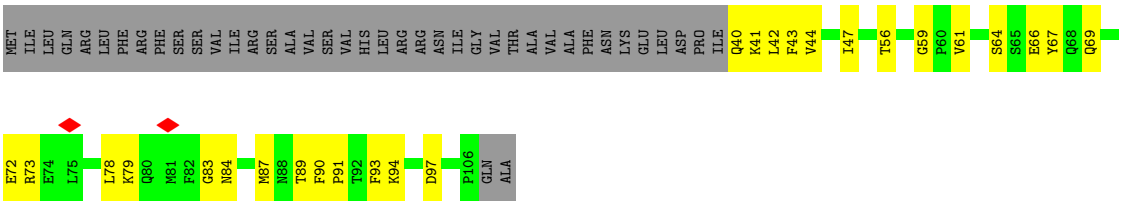




• Molecule 12: ATP synthase-coupling factor 6, mitochondrial

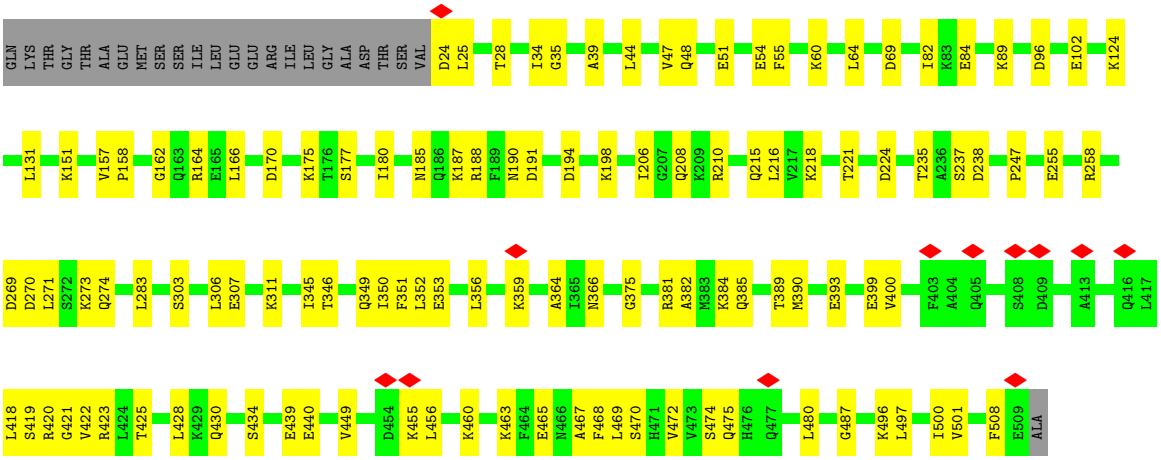
Chain L:





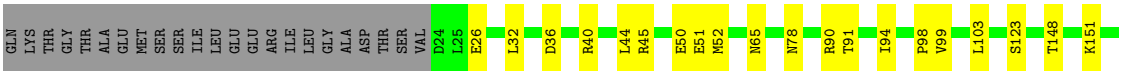
• Molecule 13: ATP synthase subunit alpha, mitochondrial

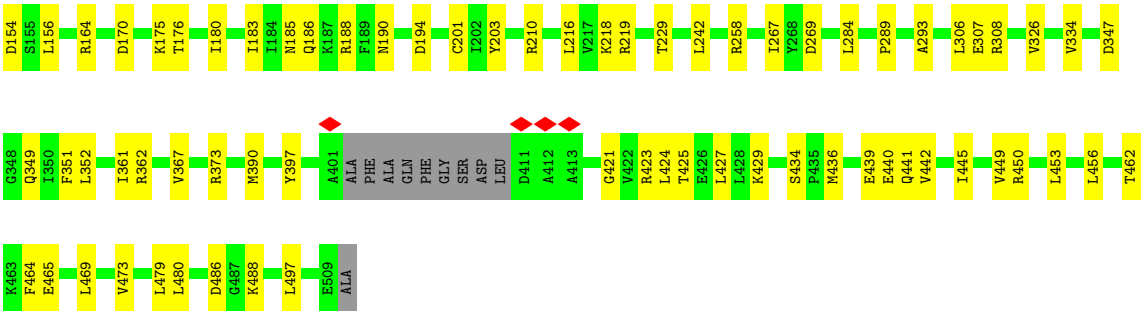
Chain A:



• Molecule 13: ATP synthase subunit alpha, mitochondrial

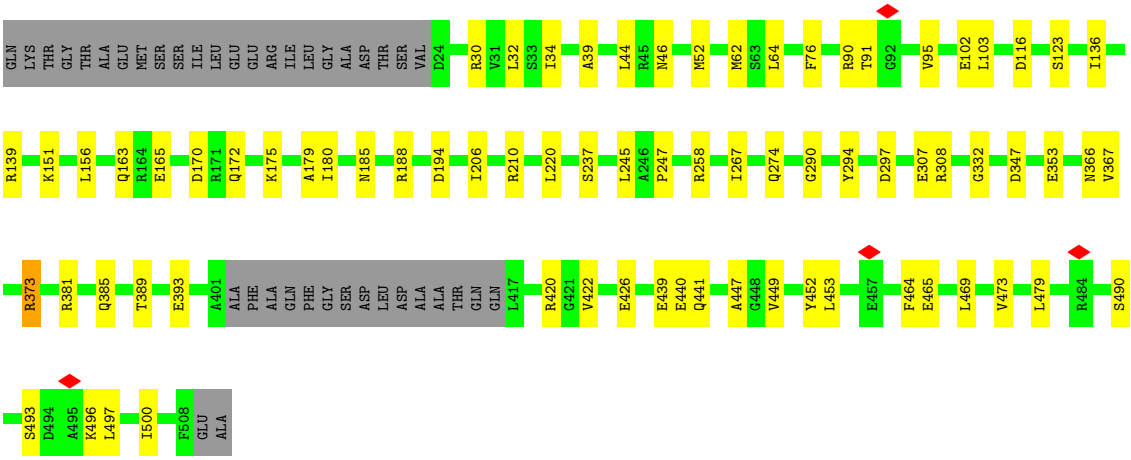
Chain B:





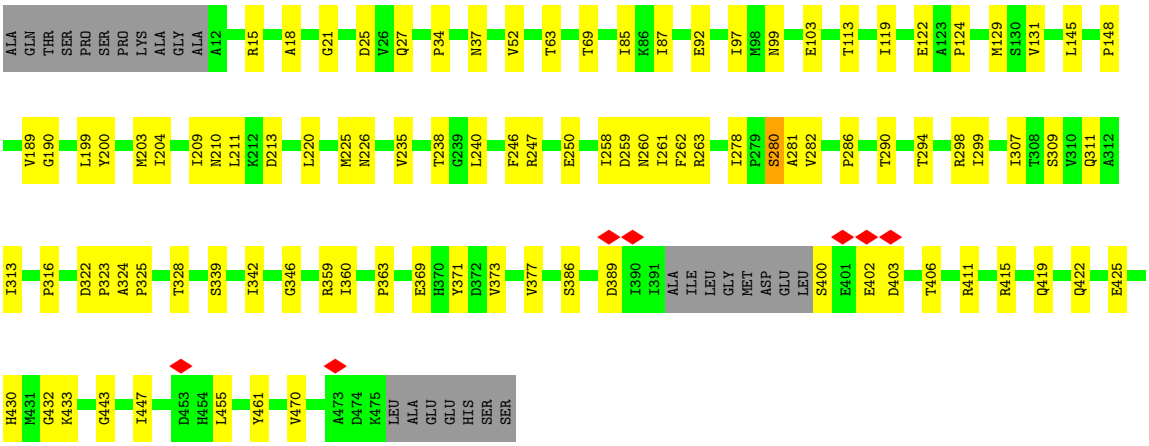
• Molecule 13: ATP synthase subunit alpha, mitochondrial

Chain C:



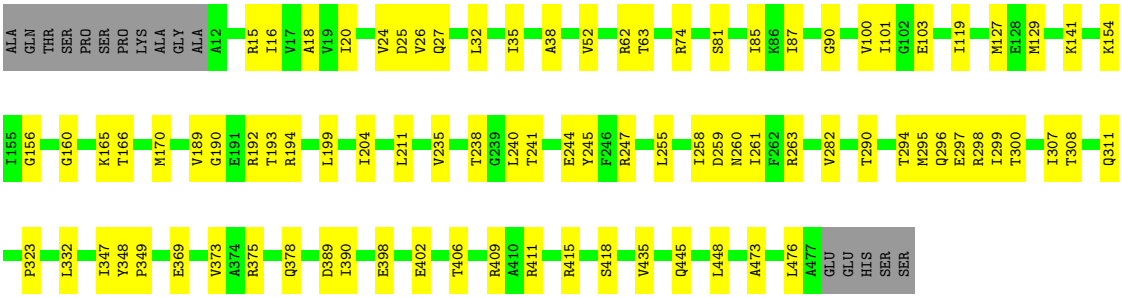
• Molecule 14: ATP synthase subunit beta, mitochondrial

Chain E:



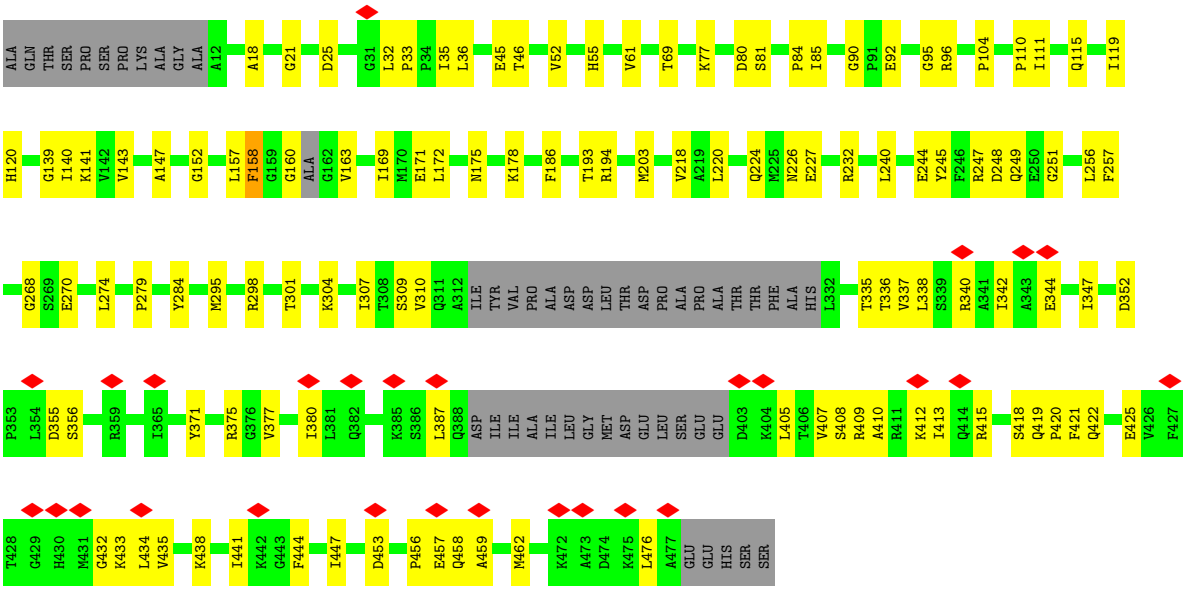
• Molecule 14: ATP synthase subunit beta, mitochondrial

Chain F:



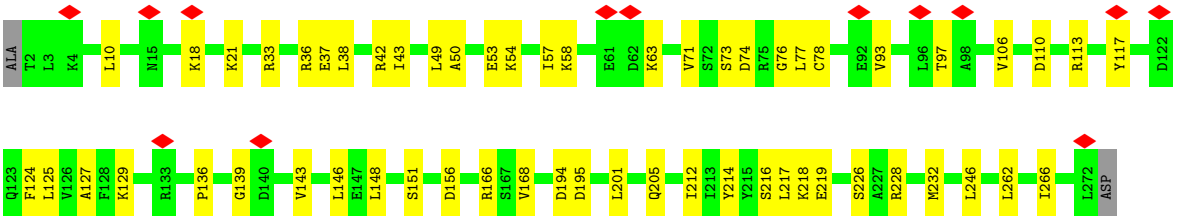
• Molecule 14: ATP synthase subunit beta, mitochondrial

Chain D:



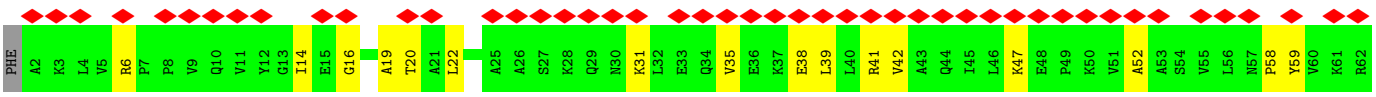
• Molecule 15: ATP synthase subunit gamma, mitochondrial

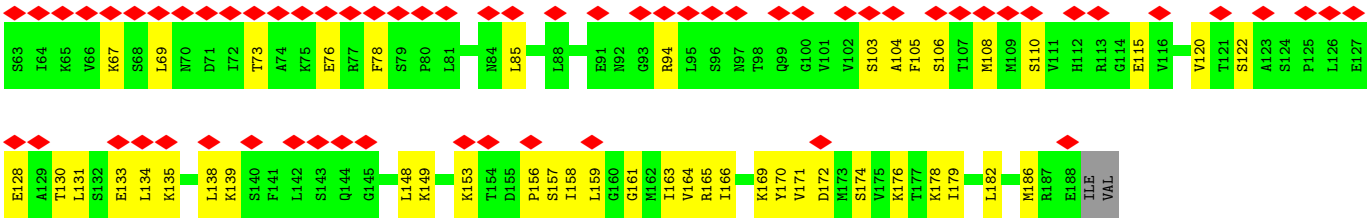
Chain G:



• Molecule 16: ATP synthase subunit O, mitochondrial

Chain O:




S63 I64 K65 V66 K67 S68 L69 N70 D71 I72 T73 A74 K75 E76 R77 F78 S79 P80 L81 N84 L85 L88 E91 N92 G93 R94 L95 S96 N97 T98 Q99 G100 V101 V102 S103 A104 F105 S106 T107 M108 M109 S110 V111 H112 R113 G114 E115 V116 V120 T121 S122 A123 S124 P125 L126 E127
E128 A129 T130 L131 S132 E133 L134 K135 L138 K139 S140 F141 L142 S143 Q144 G145 L148 K149 K153 T154 D155 P156 S157 I158 L159 G160 G161 M162 I163 V164 R165 I166 K169 Y170 V171 D172 M173 S174 V175 K176 T177 K178 I179 L182 M186 R187 E188 ILE VAL

4 Experimental information (i)

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	23272	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 ($e^-/Å^2$)	50	Depositor
Minimum defocus (nm)	1200	Depositor
Maximum defocus (nm)	2400	Depositor
Magnification	Not provided	
Image detector	FEI FALCON IV (4k x 4k)	Depositor
Maximum map value	31.451	Depositor
Minimum map value	-21.179	Depositor
Average map value	0.009	Depositor
Map value standard deviation	1.076	Depositor
Recommended contour level	4.5	Depositor
Map size (Å)	373.76, 373.76, 373.76	wwPDB
Map dimensions	512, 512, 512	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.73, 0.73, 0.73	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MG, ADP, ATP

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	1	0.43	0/543	0.55	0/732
1	2	0.41	0/543	0.54	0/732
1	3	0.42	0/543	0.54	0/732
1	4	0.42	0/543	0.53	0/732
1	5	0.42	0/543	0.56	0/732
1	6	0.41	0/543	0.55	0/732
1	7	0.41	0/543	0.54	0/732
1	8	0.42	0/543	0.56	0/732
2	H	0.37	0/987	0.42	0/1344
3	I	0.31	0/359	0.31	0/482
4	K	0.34	0/1605	0.45	0/2169
5	M	0.33	0/1291	0.41	0/1755
6	N	0.47	0/1755	0.49	0/2403
7	P	0.29	0/354	0.37	0/478
8	Q	0.55	0/437	0.63	0/598
9	R	0.37	0/640	0.42	0/858
10	S	0.21	0/619	0.38	0/841
11	T	0.20	0/354	0.35	0/480
12	L	0.16	0/560	0.35	0/748
13	A	0.34	0/3757	0.39	0/5069
13	B	0.39	0/3687	0.39	0/4973
13	C	0.39	0/3635	0.41	0/4902
14	D	0.33	0/3321	0.41	0/4494
14	E	0.37	0/3515	0.37	0/4766
14	F	0.40	0/3587	0.40	0/4865
15	G	0.33	0/2129	0.40	0/2860
16	O	0.18	0/1453	0.37	0/1958
All	All	0.37	0/38389	0.42	0/51899

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

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	1	534	0	553	10	0
1	2	534	0	553	9	0
1	3	534	0	553	14	0
1	4	534	0	553	27	0
1	5	534	0	553	57	0
1	6	534	0	553	44	0
1	7	534	0	553	13	0
1	8	534	0	553	10	0
2	H	975	0	979	28	0
3	I	354	0	364	10	0
4	K	1573	0	1552	67	0
5	M	1259	0	1238	53	0
6	N	1718	0	1879	58	0
7	P	344	0	361	6	0
8	Q	422	0	450	22	0
9	R	621	0	618	18	0
10	S	605	0	631	28	0
11	T	346	0	355	10	0
12	L	549	0	533	22	0
13	A	3706	0	3802	80	0
13	B	3639	0	3742	59	0
13	C	3587	0	3699	55	0
14	D	3272	0	3333	87	0
14	E	3458	0	3508	61	0
14	F	3529	0	3585	55	0
15	G	2103	0	2174	40	0
16	O	1437	0	1536	49	0
17	A	31	0	12	1	0
17	B	31	0	12	3	0
17	C	31	0	12	1	0
18	A	1	0	0	0	0
18	B	1	0	0	0	0
18	C	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
18	D	1	0	0	0	0
18	F	1	0	0	0	0
19	D	27	0	12	3	0
19	F	27	0	12	1	0
All	All	37921	0	38823	828	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 11.

All (828) 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)
14:D:141:LYS:HG2	14:D:435:VAL:HG21	1.56	0.87
2:H:40:THR:HG22	2:H:42:THR:H	1.46	0.81
4:K:95:GLU:O	4:K:99:ALA:HB2	1.81	0.81
14:F:170:MET:HE1	14:F:199:LEU:HD13	1.63	0.81
14:D:247:ARG:HD3	14:D:307:ILE:HG13	1.61	0.81
1:5:62:LEU:HD11	6:N:151:ILE:HD13	1.64	0.80
1:2:56:LEU:O	1:2:59:ALA:HB3	1.82	0.80
1:1:56:LEU:O	1:1:59:ALA:HB3	1.82	0.79
1:6:56:LEU:O	1:6:59:ALA:HB3	1.82	0.79
1:3:56:LEU:O	1:3:59:ALA:HB3	1.82	0.79
5:M:108:LYS:HA	5:M:111:ILE:HD12	1.64	0.79
1:8:56:LEU:O	1:8:59:ALA:HB3	1.82	0.79
1:4:56:LEU:O	1:4:59:ALA:HB3	1.82	0.79
1:5:56:LEU:O	1:5:59:ALA:HB3	1.82	0.79
1:7:56:LEU:O	1:7:59:ALA:HB3	1.82	0.78
1:7:37:ALA:HB2	1:8:46:LEU:HD21	1.65	0.78
5:M:107:SER:HA	5:M:110:ARG:HD3	1.64	0.78
1:4:37:ALA:HB2	1:5:46:LEU:HD21	1.66	0.77
14:E:260:ASN:HB3	14:E:263:ARG:HG3	1.66	0.77
1:5:34:ILE:O	1:5:38:ARG:HG3	1.85	0.76
5:M:130:MET:HG3	5:M:134:ASP:HB2	1.68	0.75
3:I:24:ARG:HA	3:I:27:LEU:HD12	1.68	0.75
1:5:38:ARG:HG2	1:6:38:ARG:NH2	2.02	0.75
2:H:122:THR:O	2:H:128:ARG:NH2	2.20	0.74
2:H:134:ARG:O	2:H:138:ASN:ND2	2.21	0.74
14:D:408:SER:OG	14:D:412:LYS:NZ	2.22	0.73
15:G:78:CYS:SG	15:G:228:ARG:NH2	2.61	0.73
13:A:420:ARG:NH1	13:A:449:VAL:O	2.22	0.72
6:N:106:ILE:O	6:N:110:ALA:HB2	1.90	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
14:D:152:GLY:HA3	14:D:301:THR:HG23	1.72	0.72
14:D:158:PHE:CE2	14:D:335:THR:HG23	2.25	0.71
8:Q:39:PRO:O	8:Q:41:PRO:HD3	1.89	0.71
16:O:14:ILE:HD13	16:O:94:ARG:HH11	1.56	0.71
6:N:31:ILE:HA	6:N:46:GLN:HE22	1.56	0.70
14:F:18:ALA:HB3	14:F:25:ASP:HB2	1.71	0.70
5:M:55:TYR:O	5:M:59:ASN:ND2	2.25	0.70
4:K:103:HIS:HE1	8:Q:43:LYS:H	1.40	0.70
13:A:187:LYS:NZ	13:A:191:ASP:OD2	2.25	0.70
1:5:38:ARG:HG2	1:6:38:ARG:HH21	1.57	0.69
4:K:4:PRO:HB3	5:M:129:GLN:HE21	1.57	0.69
6:N:40:ASN:HD22	9:R:56:ASN:HA	1.57	0.69
13:A:381:ARG:HG3	13:A:385:GLN:HE21	1.57	0.69
13:B:436:MET:O	13:B:441:GLN:NE2	2.24	0.69
13:C:95:VAL:HG11	13:C:245:LEU:HD21	1.75	0.69
14:E:260:ASN:OD1	14:E:261:ILE:N	2.25	0.69
13:C:353:GLU:OE2	13:C:366:ASN:ND2	2.25	0.69
4:K:12:VAL:HA	4:K:18:PRO:HA	1.75	0.69
6:N:41:ARG:NH1	9:R:60:GLY:O	2.26	0.69
14:D:456:PRO:HG2	14:D:459:ALA:HB2	1.76	0.68
12:L:42:LEU:HD23	16:O:182:LEU:HD21	1.74	0.68
13:A:170:ASP:O	13:A:175:LYS:NZ	2.26	0.68
13:B:185:ASN:OD1	13:B:188:ARG:NH1	2.26	0.68
14:E:419:GLN:NE2	14:E:433:LYS:O	2.25	0.68
1:4:16:VAL:HG22	1:5:13:ALA:O	1.94	0.67
4:K:96:ALA:HA	8:Q:39:PRO:HG2	1.76	0.67
12:L:42:LEU:HD21	16:O:178:LYS:HD2	1.76	0.67
13:A:175:LYS:HG2	13:A:352:LEU:HD12	1.75	0.67
13:B:170:ASP:O	13:B:175:LYS:NZ	2.27	0.67
13:A:456:LEU:HD11	13:A:460:LYS:HD2	1.77	0.67
13:B:175:LYS:HG2	13:B:352:LEU:HD12	1.75	0.67
14:F:52:VAL:HA	14:F:63:THR:HG22	1.77	0.67
14:D:80:ASP:OD2	14:D:81:SER:N	2.27	0.67
7:P:4:SER:HA	7:P:7:LYS:HB3	1.77	0.66
4:K:103:HIS:HE1	8:Q:43:LYS:N	1.92	0.66
1:5:5:ALA:HA	1:6:6:ALA:CB	2.25	0.66
5:M:18:ILE:HG23	8:Q:51:TRP:CZ2	2.30	0.66
14:E:247:ARG:HD3	14:E:307:ILE:HG13	1.77	0.66
2:H:64:VAL:HG22	2:H:74:LYS:HG2	1.77	0.66
6:N:58:MET:HE1	6:N:72:LEU:HD12	1.78	0.66
14:E:199:LEU:HD11	14:E:203:MET:HE3	1.78	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
14:D:141:LYS:CG	14:D:435:VAL:HG21	2.26	0.66
13:C:210:ARG:NH2	13:C:237:SER:OG	2.27	0.66
1:5:54:PHE:CE1	1:6:56:LEU:HD13	2.32	0.65
14:E:15:ARG:NH2	14:E:27:GLN:OE1	2.30	0.65
1:8:38:ARG:HA	1:1:38:ARG:HE	1.62	0.65
14:E:190:GLY:O	14:E:263:ARG:NH1	2.30	0.65
16:O:130:THR:HA	16:O:133:GLU:HG2	1.78	0.65
2:H:58:LEU:HD11	2:H:92:LEU:HD11	1.79	0.65
13:C:185:ASN:OD1	13:C:188:ARG:NH1	2.28	0.65
14:D:186:PHE:HB3	14:D:220:LEU:HD23	1.79	0.65
1:4:20:GLY:HA3	1:3:19:ALA:HA	1.79	0.65
9:R:24:GLU:HB3	10:S:68:THR:HA	1.78	0.64
13:A:349:GLN:HE21	13:A:351:PHE:HB2	1.63	0.64
16:O:161:GLY:HA3	16:O:174:SER:HA	1.80	0.64
4:K:158:HIS:NE2	12:L:87:MET:O	2.31	0.64
13:B:258:ARG:NH1	13:B:308:ARG:O	2.31	0.64
13:B:186:GLN:O	13:B:190:ASN:ND2	2.28	0.64
6:N:97:GLN:O	6:N:101:ASN:ND2	2.31	0.64
6:N:106:ILE:O	6:N:110:ALA:CB	2.45	0.64
13:A:185:ASN:OD1	13:A:188:ARG:NH1	2.31	0.63
1:5:5:ALA:HA	1:6:6:ALA:HB3	1.80	0.63
6:N:184:ILE:HG22	6:N:185:ASN:H	1.63	0.63
10:S:28:THR:HA	10:S:31:TYR:HB3	1.81	0.63
6:N:83:ASN:HB3	6:N:207:ALA:HB1	1.80	0.63
6:N:143:ILE:O	6:N:147:ILE:HG12	1.98	0.63
4:K:154:ARG:NH1	12:L:89:THR:O	2.32	0.63
14:D:256:LEU:O	14:D:309:SER:HA	1.99	0.63
13:A:102:GLU:OE2	13:A:102:GLU:N	2.29	0.62
13:B:44:LEU:O	14:F:74:ARG:NH1	2.31	0.62
10:S:36:GLU:O	11:T:12:LYS:NZ	2.32	0.62
1:5:5:ALA:O	1:6:6:ALA:HB1	1.97	0.62
1:5:63:PHE:O	1:5:66:MET:HB3	1.99	0.62
1:1:63:PHE:O	1:1:66:MET:HB3	1.99	0.62
1:4:63:PHE:O	1:4:66:MET:HB3	2.00	0.62
1:5:5:ALA:HB1	1:6:6:ALA:HB2	1.80	0.62
1:5:73:PHE:CZ	1:6:75:MET:HE3	2.35	0.62
9:R:61:SER:OG	9:R:62:ILE:N	2.32	0.62
14:F:15:ARG:NH2	14:F:27:GLN:OE1	2.33	0.62
14:D:18:ALA:HB3	14:D:25:ASP:HB2	1.80	0.62
16:O:164:VAL:HB	16:O:171:VAL:HG22	1.80	0.62
1:3:63:PHE:O	1:3:66:MET:HB3	1.99	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:109:LYS:HZ1	2:H:113:GLU:HB2	1.65	0.62
10:S:53:LYS:HA	10:S:56:VAL:HG12	1.82	0.62
1:7:63:PHE:O	1:7:66:MET:HB3	1.99	0.62
1:5:33:ILE:HG21	1:6:32:LEU:HA	1.81	0.62
13:C:46:ASN:O	13:C:90:ARG:NH1	2.33	0.62
16:O:158:ILE:HA	16:O:176:LYS:HG3	1.82	0.62
13:B:151:LYS:NZ	13:B:427:LEU:O	2.28	0.61
16:O:22:LEU:HD22	16:O:85:LEU:HD22	1.80	0.61
1:4:34:ILE:HG23	1:5:38:ARG:NH2	2.14	0.61
13:B:210:ARG:NH1	14:E:124:PRO:O	2.34	0.61
1:6:63:PHE:O	1:6:66:MET:HB3	1.99	0.61
14:F:244:GLU:OE2	14:F:298:ARG:NH1	2.33	0.61
14:D:420:PRO:HG2	14:D:433:LYS:H	1.65	0.61
13:A:356:LEU:HA	13:A:359:LYS:HG2	1.83	0.61
1:2:63:PHE:O	1:2:66:MET:HB3	2.00	0.61
4:K:174:ILE:HD11	16:O:186:MET:HB3	1.82	0.61
15:G:77:LEU:HB3	15:G:232:MET:HE3	1.83	0.61
4:K:143:ARG:NH2	12:L:97:ASP:OD1	2.34	0.61
9:R:76:SER:HA	9:R:79:PHE:CE1	2.36	0.61
5:M:95:GLU:HA	5:M:98:LYS:HD2	1.81	0.61
14:F:156:GLY:HA3	14:F:332:LEU:HD13	1.82	0.61
1:8:63:PHE:O	1:8:66:MET:HB3	1.99	0.61
15:G:42:ARG:NH2	15:G:219:GLU:OE2	2.32	0.61
7:P:28:MET:HE2	8:Q:17:LEU:HD22	1.82	0.60
13:C:440:GLU:HB3	13:C:469:LEU:HD11	1.83	0.60
13:C:30:ARG:CZ	16:O:59:TYR:HB2	2.31	0.60
14:F:258:ILE:HG21	14:F:261:ILE:HD13	1.83	0.60
13:C:420:ARG:NH1	13:C:449:VAL:O	2.34	0.60
13:A:356:LEU:HB2	13:A:364:ALA:HB1	1.81	0.60
14:E:189:VAL:HG22	14:E:235:VAL:HG13	1.84	0.60
1:4:18:VAL:HG12	1:5:60:MET:HB3	1.83	0.59
13:C:389:THR:HG23	13:C:449:VAL:HG21	1.84	0.59
4:K:172:HIS:HA	4:K:175:ASN:HD21	1.67	0.59
1:5:19:ALA:HA	1:6:20:GLY:HA3	1.84	0.59
13:A:206:ILE:HD11	13:A:247:PRO:HG3	1.83	0.59
16:O:159:LEU:HG	16:O:176:LYS:HD3	1.84	0.59
13:C:422:VAL:O	13:C:426:GLU:HG2	2.02	0.59
1:5:1:ASP:N	1:6:3:ASP:OD1	2.36	0.59
13:A:399:GLU:HG3	13:A:400:VAL:HG13	1.85	0.59
14:F:240:LEU:HD11	14:F:299:ILE:HD11	1.84	0.59
15:G:33:ARG:HG2	15:G:36:ARG:HH21	1.67	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:K:121:ARG:HA	5:M:18:ILE:HD13	1.83	0.58
13:A:187:LYS:HE2	13:A:224:ASP:HB3	1.83	0.58
13:B:390:MET:HE3	13:B:424:LEU:HD22	1.83	0.58
4:K:98:GLN:HA	4:K:101:ILE:HD12	1.85	0.58
13:A:235:THR:N	13:A:238:ASP:OD2	2.35	0.58
1:7:70:LEU:CD1	6:N:173:LEU:HD13	2.33	0.58
15:G:168:VAL:O	15:G:226:SER:OG	2.21	0.58
1:1:65:LEU:O	1:1:68:ALA:HB3	2.04	0.58
1:3:65:LEU:O	1:3:68:ALA:HB3	2.04	0.58
14:D:35:ILE:HG22	14:D:36:LEU:HG	1.85	0.58
1:8:65:LEU:O	1:8:68:ALA:HB3	2.04	0.58
4:K:53:ILE:HA	4:K:57:THR:HG21	1.86	0.58
13:A:208:GLN:NE2	13:A:269:ASP:O	2.37	0.58
13:B:289:PRO:HB2	13:B:293:ALA:HA	1.86	0.58
14:D:377:VAL:HA	14:D:380:ILE:HG22	1.84	0.58
15:G:49:LEU:O	15:G:53:GLU:HG2	2.04	0.58
13:C:179:ALA:HB1	13:C:267:ILE:HD13	1.84	0.57
14:D:337:VAL:HG23	14:D:352:ASP:HB3	1.85	0.57
1:7:65:LEU:O	1:7:68:ALA:HB3	2.04	0.57
1:2:65:LEU:O	1:2:68:ALA:HB3	2.04	0.57
14:D:140:ILE:HA	14:D:419:GLN:HE22	1.69	0.57
2:H:127:THR:O	2:H:131:ILE:HG12	2.04	0.57
15:G:76:GLY:O	15:G:228:ARG:NH2	2.26	0.57
13:A:421:GLY:O	13:A:425:THR:HG23	2.04	0.57
14:D:420:PRO:HB2	14:D:432:GLY:HA2	1.86	0.57
1:6:65:LEU:O	1:6:68:ALA:HB3	2.04	0.57
5:M:53:TRP:O	5:M:57:LYS:N	2.35	0.57
15:G:106:VAL:HG21	15:G:148:LEU:HD11	1.86	0.57
14:E:278:ILE:HG23	15:G:266:ILE:HD12	1.86	0.57
14:D:203:MET:HE1	14:D:220:LEU:HD21	1.87	0.57
14:D:240:LEU:HD21	14:D:298:ARG:HB2	1.87	0.57
4:K:91:ALA:O	4:K:95:GLU:HG2	2.04	0.57
13:C:172:GLN:NE2	17:C:601:ATP:O3G	2.38	0.57
13:B:78:ASN:ND2	14:E:122:GLU:OE2	2.37	0.57
13:B:479:LEU:HD21	13:B:497:LEU:HD23	1.87	0.57
14:E:203:MET:HB3	14:E:209:ILE:HG13	1.85	0.57
1:4:65:LEU:O	1:4:68:ALA:HB3	2.04	0.57
1:5:65:LEU:O	1:5:68:ALA:HB3	2.04	0.57
15:G:113:ARG:O	15:G:117:TYR:HB2	2.05	0.57
16:O:179:ILE:HA	16:O:182:LEU:HG	1.87	0.57
5:M:57:LYS:HZ2	5:M:66:VAL:HG13	1.70	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:N:112:THR:HG22	6:N:225:ASN:HD22	1.70	0.56
13:B:421:GLY:O	13:B:425:THR:HG23	2.05	0.56
14:D:268:GLY:HA3	14:D:284:TYR:CE2	2.41	0.56
2:H:132:GLN:HA	2:H:135:ILE:HD12	1.87	0.56
13:B:180:ILE:HD11	13:B:216:LEU:HD11	1.86	0.56
16:O:128:GLU:HA	16:O:131:LEU:HB3	1.86	0.56
13:C:381:ARG:HG3	13:C:385:GLN:HE21	1.70	0.56
6:N:140:MET:SD	6:N:140:MET:N	2.79	0.56
1:7:70:LEU:HD11	6:N:173:LEU:HD13	1.87	0.56
10:S:85:TRP:CD1	11:T:15:ARG:HG2	2.40	0.56
13:C:479:LEU:HD21	13:C:497:LEU:HB3	1.87	0.56
14:E:455:LEU:HD11	14:E:470:VAL:HG12	1.86	0.56
14:D:444:PHE:HA	14:D:447:ILE:HG22	1.88	0.56
13:C:297:ASP:HB3	14:D:274:LEU:HD21	1.88	0.56
1:4:33:ILE:HD11	1:5:50:ALA:N	2.22	0.55
14:F:25:ASP:OD2	14:F:62:ARG:NH2	2.38	0.55
5:M:21:GLN:HA	5:M:24:LYS:HD3	1.88	0.55
1:5:72:LEU:CD2	1:6:71:ILE:HD11	2.37	0.55
2:H:138:ASN:HA	2:H:141:LEU:HD12	1.88	0.55
13:A:28:THR:HG22	13:A:89:LYS:HG3	1.87	0.55
14:D:45:GLU:HG3	14:D:46:THR:HG23	1.88	0.55
6:N:99:SER:HG	8:Q:9:TRP:CG	2.24	0.55
11:T:17:SER:O	11:T:21:LEU:HG	2.07	0.55
14:D:420:PRO:HB3	14:D:462:MET:HE1	1.89	0.55
13:B:423:ARG:NH2	13:B:456:LEU:O	2.30	0.55
14:D:377:VAL:HG23	14:D:413:ILE:HD12	1.87	0.55
13:A:215:GLN:O	13:A:218:LYS:HG3	2.07	0.55
13:C:62:MET:HB2	13:C:76:PHE:HE2	1.72	0.55
14:F:369:GLU:OE2	14:F:445:GLN:NE2	2.35	0.55
1:5:47:PHE:HD1	1:6:49:TYR:CZ	2.25	0.55
15:G:125:LEU:HD21	15:G:151:SER:HB2	1.88	0.55
15:G:194:ASP:OD1	15:G:195:ASP:N	2.40	0.55
6:N:142:VAL:O	6:N:146:THR:HG23	2.07	0.55
13:C:496:LYS:O	13:C:500:ILE:HG12	2.07	0.55
11:T:19:LEU:O	11:T:23:VAL:HG23	2.06	0.54
14:D:147:ALA:HB2	14:D:356:SER:HB3	1.87	0.54
13:B:99:VAL:O	13:B:123:SER:OG	2.24	0.54
13:B:486:ASP:HB3	13:B:488:LYS:HG2	1.89	0.54
13:C:52:MET:O	13:C:91:THR:OG1	2.25	0.54
14:D:405:LEU:O	14:D:409:ARG:HG2	2.08	0.54
1:3:1:ASP:OD1	1:3:2:ILE:N	2.41	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:M:128:ASP:OD1	5:M:129:GLN:NE2	2.41	0.54
14:E:52:VAL:HA	14:E:63:THR:HG22	1.87	0.54
1:5:1:ASP:OD1	1:5:2:ILE:N	2.41	0.54
13:A:84:GLU:OE2	14:D:55:HIS:ND1	2.31	0.54
14:D:244:GLU:OE2	14:D:298:ARG:HD2	2.08	0.54
2:H:18:PHE:CE2	2:H:20:PHE:HB2	2.43	0.54
4:K:162:GLN:NE2	5:M:59:ASN:O	2.33	0.54
14:D:84:PRO:O	14:D:85:ILE:HG13	2.07	0.54
1:6:1:ASP:OD1	1:6:2:ILE:N	2.41	0.54
16:O:120:VAL:HG21	16:O:134:LEU:HD21	1.88	0.54
1:4:1:ASP:OD1	1:4:2:ILE:N	2.41	0.54
14:E:280:SER:OG	14:E:281:ALA:N	2.40	0.54
1:5:62:LEU:CD1	6:N:151:ILE:HD13	2.36	0.54
13:A:467:ALA:O	13:A:470:SER:OG	2.24	0.54
13:B:36:ASP:HB3	13:B:284:LEU:HB3	1.89	0.54
14:F:141:LYS:HD3	14:F:435:VAL:HG21	1.89	0.54
5:M:53:TRP:NE1	5:M:70:GLU:OE1	2.41	0.54
16:O:134:LEU:O	16:O:138:LEU:HG	2.08	0.54
1:1:1:ASP:OD1	1:1:2:ILE:N	2.41	0.53
13:A:210:ARG:NH1	13:A:237:SER:OG	2.41	0.53
10:S:46:PRO:O	10:S:49:ILE:HG12	2.08	0.53
10:S:102:VAL:HG11	11:T:34:LYS:HD3	1.89	0.53
1:5:65:LEU:HD11	1:6:63:PHE:CD2	2.44	0.53
1:8:1:ASP:OD1	1:8:2:ILE:N	2.41	0.53
1:2:1:ASP:OD1	1:2:2:ILE:N	2.41	0.53
11:T:38:GLU:HA	11:T:41:ARG:HD2	1.90	0.53
14:E:322:ASP:HB3	14:E:325:PRO:HD2	1.89	0.53
1:7:1:ASP:OD1	1:7:2:ILE:N	2.41	0.53
2:H:40:THR:HB	2:H:43:GLY:O	2.09	0.53
10:S:86:PHE:HB2	11:T:18:ALA:HB1	1.91	0.53
13:C:116:ASP:HA	14:F:127:MET:HE3	1.89	0.53
13:C:453:LEU:HD21	13:C:464:PHE:CE2	2.44	0.53
14:E:204:ILE:HD13	14:E:211:LEU:HD11	1.91	0.53
14:F:38:ALA:O	14:F:81:SER:OG	2.24	0.53
13:C:170:ASP:OD2	13:C:332:GLY:N	2.42	0.53
14:E:240:LEU:HD21	14:E:298:ARG:HB2	1.90	0.53
14:F:100:VAL:HG13	14:F:101:ILE:HG23	1.91	0.53
14:D:21:GLY:O	14:D:69:THR:OG1	2.23	0.53
14:F:411:ARG:O	14:F:415:ARG:HG2	2.09	0.53
13:A:177:SER:OG	17:A:601:ATP:O1A	2.23	0.52
13:B:445:ILE:O	13:B:449:VAL:HG22	2.10	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
13:C:490:SER:O	13:C:493:SER:N	2.41	0.52
14:D:35:ILE:HD12	14:D:35:ILE:H	1.74	0.52
14:D:257:PHE:CD1	14:D:310:VAL:HB	2.45	0.52
6:N:103:ALA:HB2	8:Q:13:ILE:HD12	1.90	0.52
14:D:158:PHE:CD2	14:D:335:THR:HG23	2.43	0.52
15:G:73:SER:OG	15:G:74:ASP:N	2.41	0.52
5:M:34:ASN:O	5:M:38:THR:HG23	2.08	0.52
14:F:296:GLN:HE22	14:F:311:GLN:HE22	1.58	0.52
6:N:40:ASN:ND2	9:R:56:ASN:HA	2.24	0.52
14:E:261:ILE:HG21	14:E:313:ILE:HD13	1.91	0.52
14:E:309:SER:OG	14:E:311:GLN:NE2	2.42	0.52
14:F:160:GLY:O	14:F:165:LYS:NZ	2.42	0.52
14:F:290:THR:O	14:F:294:THR:HG23	2.10	0.52
5:M:145:ASP:OD1	5:M:148:LYS:N	2.30	0.52
6:N:123:ASN:HA	6:N:126:ALA:HB3	1.92	0.52
4:K:111:GLU:OE2	5:M:110:ARG:NH2	2.42	0.52
13:B:51:GLU:OE2	13:B:90:ARG:NH1	2.41	0.52
14:D:340:ARG:O	14:D:344:GLU:HG2	2.09	0.52
15:G:18:LYS:HA	15:G:21:LYS:HD3	1.91	0.52
12:L:42:LEU:HD12	12:L:42:LEU:H	1.73	0.52
16:O:31:LYS:O	16:O:35:VAL:HG22	2.10	0.52
5:M:27:ALA:O	5:M:30:LEU:HB3	2.09	0.52
1:5:12:GLY:HA3	1:6:13:ALA:HB3	1.91	0.52
14:F:154:LYS:NZ	14:F:296:GLN:O	2.43	0.52
1:8:21:SER:O	1:8:25:ILE:HG13	2.10	0.52
14:F:119:ILE:O	14:F:298:ARG:NH2	2.43	0.52
1:6:36:TYR:OH	1:7:45:GLN:HG2	2.10	0.51
6:N:27:PRO:HB2	6:N:28:PRO:HD3	1.92	0.51
10:S:78:VAL:O	10:S:81:GLU:HB3	2.10	0.51
14:D:387:LEU:HD13	14:D:407:VAL:HG23	1.92	0.51
1:7:21:SER:O	1:7:25:ILE:HG13	2.10	0.51
1:2:21:SER:O	1:2:25:ILE:HG13	2.10	0.51
5:M:103:TRP:CD2	8:Q:50:PRO:HG3	2.45	0.51
7:P:35:VAL:O	7:P:39:ARG:HG3	2.10	0.51
1:3:40:PRO:HD2	3:I:2:VAL:HB	1.92	0.51
4:K:155:LEU:HD22	5:M:56:TYR:CD1	2.45	0.51
9:R:48:TYR:O	9:R:52:ASN:ND2	2.44	0.51
14:D:456:PRO:HG3	14:D:476:LEU:HD22	1.92	0.51
1:4:21:SER:O	1:4:25:ILE:HG13	2.10	0.51
13:C:258:ARG:NH1	13:C:308:ARG:O	2.44	0.51
1:5:5:ALA:CA	1:6:6:ALA:CB	2.88	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:R:31:MET:SD	9:R:32:ARG:NH2	2.84	0.51
13:C:32:LEU:HD21	16:O:58:PRO:HB2	1.93	0.51
1:1:21:SER:O	1:1:25:ILE:HG13	2.10	0.51
4:K:143:ARG:NH1	5:M:79:PRO:O	2.43	0.51
5:M:11:ASP:OD2	5:M:12:TRP:N	2.44	0.51
10:S:84:MET:HE3	10:S:84:MET:O	2.10	0.51
13:C:102:GLU:OE1	13:C:123:SER:HA	2.11	0.51
1:6:21:SER:O	1:6:25:ILE:HG13	2.10	0.51
6:N:160:LEU:O	6:N:164:ILE:HG12	2.11	0.51
14:D:247:ARG:NE	14:D:248:ASP:OD1	2.27	0.51
15:G:57:ILE:C	15:G:58:LYS:HD2	2.35	0.51
1:6:62:LEU:HD21	6:N:162:ALA:HB3	1.91	0.51
14:F:166:THR:O	14:F:170:MET:HG2	2.10	0.51
16:O:103:SER:O	16:O:106:SER:OG	2.27	0.51
4:K:118:VAL:HG22	8:Q:51:TRP:CZ3	2.46	0.51
10:S:67:LEU:HD11	10:S:71:GLU:HB2	1.92	0.51
13:A:440:GLU:HB3	13:A:469:LEU:HD11	1.92	0.51
1:3:21:SER:O	1:3:25:ILE:HG13	2.10	0.51
6:N:101:ASN:ND2	6:N:161:THR:OG1	2.44	0.51
13:C:206:ILE:HD11	13:C:247:PRO:HG3	1.93	0.51
14:F:129:MET:HE1	14:F:300:THR:HG21	1.93	0.51
15:G:38:LEU:O	15:G:42:ARG:N	2.32	0.51
1:5:21:SER:O	1:5:25:ILE:HG13	2.10	0.50
13:A:54:GLU:HG2	13:A:60:LYS:HD3	1.93	0.50
1:5:15:THR:HG21	1:6:14:ALA:HB1	1.92	0.50
13:A:496:LYS:O	13:A:500:ILE:HG13	2.10	0.50
4:K:140:TYR:CZ	4:K:144:LEU:HD11	2.45	0.50
6:N:5:LEU:HD11	6:N:173:LEU:HD21	1.94	0.50
13:C:44:LEU:HD22	13:C:90:ARG:HG3	1.92	0.50
10:S:88:VAL:HA	10:S:91:ILE:HG12	1.93	0.50
13:A:158:PRO:HG2	13:A:375:GLY:HA2	1.93	0.50
13:B:424:LEU:HD11	13:B:449:VAL:HG12	1.91	0.50
14:F:87:ILE:HD13	14:F:238:THR:HG23	1.92	0.50
2:H:78:SER:HB3	3:I:23:VAL:HG21	1.92	0.50
14:E:346:GLY:O	14:E:461:TYR:OH	2.28	0.50
14:F:347:ILE:HG23	14:F:418:SER:HB3	1.93	0.50
4:K:114:GLN:O	4:K:118:VAL:HG23	2.12	0.50
1:6:8:PHE:HD2	1:7:6:ALA:HB3	1.76	0.50
4:K:175:ASN:O	4:K:179:LYS:HE3	2.12	0.50
1:5:4:THR:OG1	1:6:3:ASP:OD1	2.30	0.49
6:N:110:ALA:O	6:N:114:ILE:HG12	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
13:A:393:GLU:OE1	13:A:420:ARG:NH1	2.45	0.49
14:E:369:GLU:O	14:E:373:VAL:HG23	2.12	0.49
16:O:69:LEU:O	16:O:73:THR:HG23	2.12	0.49
4:K:121:ARG:HH12	8:Q:51:TRP:CD1	2.30	0.49
4:K:172:HIS:HA	4:K:175:ASN:ND2	2.26	0.49
6:N:109:TRP:O	6:N:113:VAL:HG23	2.12	0.49
14:E:21:GLY:O	14:E:69:THR:OG1	2.27	0.49
6:N:53:THR:HG21	6:N:76:ILE:HG21	1.94	0.49
17:B:601:ATP:O2G	14:E:359:ARG:NH2	2.38	0.49
16:O:157:SER:O	16:O:176:LYS:NZ	2.43	0.49
1:5:26:GLY:HA3	1:6:24:GLY:HA2	1.94	0.49
13:C:469:LEU:O	13:C:473:VAL:HG23	2.11	0.49
4:K:113:SER:O	4:K:117:LEU:HD13	2.12	0.49
5:M:99:SER:O	5:M:102:GLU:HG3	2.13	0.49
1:4:42:LEU:HD13	1:3:36:TYR:CE2	2.47	0.49
4:K:128:GLN:O	4:K:132:ILE:HG12	2.13	0.49
6:N:13:THR:O	9:R:82:LYS:NZ	2.40	0.49
13:A:218:LYS:HA	13:A:221:THR:HG22	1.93	0.49
13:B:439:GLU:HG3	13:B:480:LEU:HB3	1.95	0.49
14:D:141:LYS:H	14:D:419:GLN:HE22	1.60	0.49
16:O:178:LYS:O	16:O:182:LEU:HG	2.12	0.49
3:I:24:ARG:O	3:I:27:LEU:HB2	2.13	0.49
8:Q:48:ASN:ND2	8:Q:49:LYS:O	2.46	0.49
13:B:52:MET:O	13:B:91:THR:OG1	2.22	0.49
13:B:203:TYR:OH	13:B:269:ASP:OD2	2.27	0.49
4:K:44:TYR:HB2	10:S:84:MET:SD	2.52	0.49
5:M:103:TRP:CE2	8:Q:50:PRO:HG3	2.48	0.49
13:A:55:PHE:CD1	13:A:82:ILE:HG12	2.48	0.48
14:E:148:PRO:HB2	14:E:360:ILE:HD11	1.95	0.48
14:F:282:VAL:HG21	14:F:323:PRO:CD	2.43	0.48
14:D:232:ARG:NH2	14:D:270:GLU:OE1	2.33	0.48
15:G:124:PHE:HZ	15:G:127:ALA:HB2	1.77	0.48
13:A:460:LYS:HD3	13:A:508:PHE:HE1	1.78	0.48
14:E:18:ALA:HB3	14:E:25:ASP:HB2	1.95	0.48
14:F:375:ARG:NH1	14:F:378:GLN:OE1	2.43	0.48
14:D:337:VAL:HG21	14:D:355:ASP:HB2	1.94	0.48
12:L:66:GLU:HG3	12:L:69:GLN:HE22	1.77	0.48
14:F:204:ILE:HD13	14:F:211:LEU:HD11	1.96	0.48
14:F:402:GLU:O	14:F:406:THR:HG23	2.14	0.48
1:5:27:THR:HA	1:6:27:THR:HG21	1.96	0.48
13:A:162:GLY:HA3	13:A:311:LYS:HB2	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
13:A:247:PRO:HG2	13:A:274:GLN:HG3	1.95	0.48
15:G:37:GLU:N	15:G:37:GLU:OE1	2.47	0.48
16:O:38:GLU:OE1	16:O:78:PHE:HA	2.14	0.48
10:S:90:GLU:HA	11:T:22:GLY:O	2.13	0.48
13:A:157:VAL:HG11	13:A:350:ILE:HG12	1.94	0.48
13:A:455:LYS:HD2	13:A:455:LYS:N	2.29	0.48
14:E:400:SER:N	14:E:403:ASP:HB2	2.29	0.48
4:K:126:ASP:HB3	4:K:129:ARG:HH21	1.79	0.48
14:D:33:PRO:HB3	14:D:80:ASP:OD1	2.14	0.48
14:D:409:ARG:CZ	14:D:453:ASP:HB2	2.44	0.48
13:B:349:GLN:HG3	13:B:351:PHE:CE2	2.49	0.48
13:C:393:GLU:OE1	13:C:420:ARG:NH1	2.46	0.48
4:K:161:VAL:HG12	12:L:78:LEU:HD23	1.95	0.48
6:N:55:LYS:O	6:N:59:THR:OG1	2.25	0.48
13:B:307:GLU:OE1	14:F:193:THR:OG1	2.20	0.48
14:E:363:PRO:HD3	14:E:371:TYR:CD1	2.49	0.48
14:D:141:LYS:H	14:D:419:GLN:NE2	2.11	0.48
14:D:457:GLU:OE1	14:D:458:GLN:NE2	2.47	0.48
1:5:47:PHE:HD1	1:6:49:TYR:HH	1.61	0.48
13:C:30:ARG:NH1	16:O:58:PRO:HG2	2.28	0.48
14:D:347:ILE:HG23	14:D:418:SER:HB2	1.96	0.48
13:A:353:GLU:CD	13:A:366:ASN:HD22	2.22	0.48
13:A:390:MET:HE1	13:A:428:LEU:HD21	1.95	0.48
5:M:152:TRP:CD1	5:M:153:PRO:HA	2.49	0.47
1:4:19:ALA:HB1	1:5:20:GLY:HA3	1.95	0.47
1:5:62:LEU:HD22	6:N:151:ILE:HG21	1.96	0.47
2:H:138:ASN:O	2:H:142:VAL:HG23	2.14	0.47
14:E:299:ILE:HG23	14:E:307:ILE:HG22	1.96	0.47
16:O:156:PRO:O	16:O:159:LEU:HD23	2.14	0.47
1:5:72:LEU:HD22	1:6:71:ILE:HD11	1.96	0.47
6:N:140:MET:O	6:N:144:ILE:HG22	2.14	0.47
13:B:361:ILE:HG13	13:B:429:LYS:HE2	1.96	0.47
1:6:62:LEU:HD11	6:N:162:ALA:C	2.39	0.47
4:K:147:VAL:HG23	12:L:93:PHE:CZ	2.49	0.47
13:C:373:ARG:NH1	19:D:501:ADP:N7	2.62	0.47
14:E:425:GLU:HG2	14:E:430:HIS:O	2.15	0.47
14:F:389:ASP:OD1	14:F:390:ILE:HD12	2.14	0.47
14:F:409:ARG:NH1	14:F:448:LEU:O	2.47	0.47
5:M:149:TYR:CE2	5:M:156:PRO:HD3	2.49	0.47
13:B:424:LEU:HD23	13:B:424:LEU:HA	1.70	0.47
4:K:156:ASP:HA	4:K:159:ILE:HG22	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
10:S:71:GLU:N	10:S:71:GLU:OE1	2.48	0.47
14:E:34:PRO:HD2	14:E:37:ASN:ND2	2.30	0.47
1:5:54:PHE:CD1	1:6:56:LEU:HD13	2.49	0.47
4:K:130:ASN:HD21	5:M:8:LYS:H	1.62	0.47
13:A:164:ARG:NH1	13:A:307:GLU:OE1	2.47	0.47
13:A:194:ASP:N	13:A:194:ASP:OD1	2.48	0.47
13:B:98:PRO:O	13:B:103:LEU:HD11	2.15	0.47
13:B:148:THR:OG1	13:B:154:ASP:OD1	2.30	0.47
13:C:151:LYS:HG2	13:C:441:GLN:HG2	1.96	0.47
13:C:194:ASP:OD1	13:C:194:ASP:N	2.45	0.47
14:E:99:ASN:OD1	14:E:103:GLU:N	2.39	0.47
14:E:124:PRO:HB2	14:E:129:MET:HE2	1.97	0.47
14:E:325:PRO:HA	14:E:328:THR:HG22	1.95	0.47
14:E:422:GLN:HB2	14:E:432:GLY:HA3	1.96	0.47
14:F:16:ILE:HG12	14:F:26:VAL:HG12	1.97	0.47
14:D:409:ARG:NH2	14:D:453:ASP:HB2	2.30	0.47
1:2:36:TYR:OH	1:3:45:GLN:HG2	2.15	0.47
2:H:109:LYS:NZ	2:H:113:GLU:HB2	2.30	0.47
13:B:26:GLU:HG2	13:B:45:ARG:HB2	1.96	0.47
13:B:183:ILE:HD11	13:B:267:ILE:HD13	1.96	0.47
14:D:178:LYS:NZ	14:D:422:GLN:OE1	2.29	0.47
1:4:13:ALA:HB3	1:3:12:GLY:HA3	1.97	0.47
1:4:27:THR:HG21	1:3:27:THR:HA	1.97	0.47
4:K:45:ALA:O	4:K:51:TYR:N	2.35	0.47
4:K:145:TYR:CD1	5:M:48:PRO:HG3	2.49	0.47
14:E:443:GLY:O	14:E:447:ILE:HG13	2.15	0.47
14:F:398:GLU:N	14:F:398:GLU:OE1	2.48	0.47
14:D:405:LEU:HD11	14:D:409:ARG:HH21	1.79	0.47
15:G:93:VAL:O	15:G:97:THR:OG1	2.27	0.47
13:B:434:SER:O	13:B:434:SER:OG	2.25	0.47
14:D:115:GLN:N	14:D:115:GLN:OE1	2.48	0.47
1:4:18:VAL:CG1	1:5:60:MET:HE3	2.45	0.46
5:M:143:LYS:HB2	5:M:143:LYS:HE3	1.65	0.46
13:C:139:ARG:NH2	13:C:307:GLU:O	2.44	0.46
14:D:251:GLY:HA2	14:D:304:LYS:O	2.14	0.46
5:M:72:LYS:HA	5:M:75:ALA:HB3	1.97	0.46
13:A:24:ASP:OD1	13:A:24:ASP:N	2.46	0.46
16:O:110:SER:O	16:O:115:GLU:HB2	2.15	0.46
5:M:97:VAL:HA	5:M:100:CYS:SG	2.56	0.46
6:N:115:MET:HE1	6:N:226:THR:HG22	1.96	0.46
13:A:381:ARG:NH1	13:A:384:LYS:HD2	2.30	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
14:D:77:LYS:HD3	14:D:77:LYS:HA	1.78	0.46
1:4:19:ALA:HB1	1:5:20:GLY:CA	2.46	0.46
1:5:36:TYR:CZ	1:6:42:LEU:HD13	2.51	0.46
5:M:104:VAL:HA	5:M:107:SER:OG	2.15	0.46
5:M:124:LEU:HD23	5:M:124:LEU:HA	1.69	0.46
13:C:439:GLU:CD	13:C:439:GLU:H	2.24	0.46
15:G:217:LEU:HD23	15:G:217:LEU:HA	1.81	0.46
1:5:18:VAL:HB	1:6:64:CYS:SG	2.56	0.46
4:K:40:GLY:O	10:S:84:MET:HE2	2.16	0.46
10:S:50:GLN:HA	10:S:53:LYS:HG2	1.97	0.46
13:B:218:LYS:HD2	14:E:131:VAL:HB	1.97	0.46
14:E:258:ILE:HB	14:E:311:GLN:HG2	1.98	0.46
15:G:71:VAL:HG21	15:G:216:SER:OG	2.16	0.46
1:4:59:ALA:O	1:4:62:LEU:HB3	2.16	0.46
1:4:65:LEU:HD11	1:5:63:PHE:CD2	2.51	0.46
13:B:453:LEU:HD11	13:B:464:PHE:CE1	2.51	0.46
14:F:90:GLY:HA2	14:F:245:TYR:CE2	2.50	0.46
14:F:247:ARG:HD3	14:F:307:ILE:HG13	1.97	0.46
14:F:369:GLU:O	14:F:373:VAL:HG23	2.15	0.46
16:O:47:LYS:HA	16:O:52:ALA:HB2	1.97	0.46
6:N:71:MET:HE2	6:N:108:LEU:HD21	1.98	0.46
3:I:17:GLN:NE2	15:G:146:LEU:HD11	2.31	0.46
10:S:62:GLY:HA2	10:S:64:PHE:CE1	2.51	0.46
12:L:40:GLN:HB2	12:L:43:PHE:CD2	2.50	0.46
13:A:166:LEU:HB2	13:A:346:THR:HG21	1.97	0.46
13:B:397:TYR:CG	13:B:421:GLY:HA3	2.51	0.46
1:4:33:ILE:HD11	1:5:50:ALA:CA	2.46	0.46
4:K:120:LYS:HB3	5:M:18:ILE:HD11	1.98	0.46
12:L:41:LYS:HA	12:L:44:VAL:HG12	1.98	0.46
13:C:290:GLY:N	13:C:294:TYR:O	2.42	0.46
15:G:201:LEU:HD23	15:G:201:LEU:HA	1.69	0.46
4:K:121:ARG:HA	5:M:18:ILE:HG21	1.97	0.45
14:D:224:GLN:N	14:D:227:GLU:OE1	2.49	0.45
14:D:409:ARG:NH1	14:D:453:ASP:HB2	2.31	0.45
4:K:173:MET:HE2	4:K:173:MET:HA	1.97	0.45
13:B:164:ARG:NH1	13:B:306:LEU:O	2.44	0.45
14:E:280:SER:HB2	14:E:286:PRO:HA	1.98	0.45
1:4:9:ILE:CG1	1:5:6:ALA:HB1	2.46	0.45
1:1:59:ALA:O	1:1:62:LEU:HB3	2.16	0.45
4:K:13:ARG:N	4:K:17:ILE:O	2.32	0.45
13:A:89:LYS:HE2	13:A:89:LYS:HB2	1.82	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
13:B:373:ARG:HH22	14:F:192:ARG:NH1	2.13	0.45
13:C:373:ARG:HB3	19:D:501:ADP:HN62	1.82	0.45
4:K:96:ALA:HA	8:Q:39:PRO:CG	2.45	0.45
1:2:59:ALA:O	1:2:62:LEU:HB3	2.16	0.45
4:K:97:LYS:HE3	5:M:124:LEU:HB2	1.99	0.45
14:D:434:LEU:H	14:D:434:LEU:HD23	1.82	0.45
1:5:2:ILE:HA	1:6:2:ILE:HD13	1.97	0.45
4:K:39:THR:O	4:K:43:LEU:HD23	2.16	0.45
10:S:47:ARG:HA	10:S:50:GLN:OE1	2.17	0.45
13:C:103:LEU:HD23	13:C:103:LEU:HA	1.82	0.45
14:E:386:SER:O	14:E:389:ASP:HB2	2.16	0.45
1:6:8:PHE:HD2	1:7:6:ALA:CB	2.29	0.45
14:E:200:TYR:CZ	14:E:204:ILE:HD11	2.52	0.45
14:F:255:LEU:HD23	14:F:308:THR:HB	1.97	0.45
14:F:259:ASP:HA	14:F:260:ASN:HA	1.69	0.45
14:D:410:ALA:HA	14:D:413:ILE:HG12	1.99	0.45
2:H:141:LEU:O	2:H:145:LEU:HG	2.17	0.45
4:K:130:ASN:HD21	5:M:8:LYS:N	2.15	0.45
6:N:14:ILE:HG22	6:N:19:ALA:HB2	1.98	0.45
16:O:105:PHE:HA	16:O:108:MET:HG3	1.99	0.45
1:7:59:ALA:O	1:7:62:LEU:HB3	2.16	0.45
13:A:34:ILE:HG13	13:A:35:GLY:N	2.31	0.45
13:A:96:ASP:N	13:A:96:ASP:OD1	2.49	0.45
13:B:65:ASN:ND2	14:F:20:ILE:HG23	2.32	0.45
15:G:49:LEU:HD21	15:G:212:ILE:HD12	1.97	0.45
13:C:163:GLN:NE2	13:C:165:GLU:OE2	2.50	0.45
14:D:96:ARG:HD3	14:D:111:ILE:HG12	1.98	0.45
15:G:50:ALA:HB1	15:G:54:LYS:HZ2	1.82	0.45
1:5:47:PHE:CD1	1:6:49:TYR:CZ	3.04	0.44
1:3:59:ALA:O	1:3:62:LEU:HB3	2.16	0.44
2:H:23:PRO:O	15:G:214:TYR:OH	2.24	0.44
13:A:497:LEU:O	13:A:501:VAL:HG23	2.17	0.44
4:K:127:VAL:HG22	5:M:10:ILE:HG21	1.98	0.44
4:K:171:GLU:O	4:K:174:ILE:HG22	2.16	0.44
6:N:93:THR:HB	6:N:96:THR:HG23	2.00	0.44
9:R:20:VAL:O	9:R:21:LYS:HB2	2.17	0.44
14:F:189:VAL:HG22	14:F:235:VAL:HG13	1.99	0.44
16:O:104:ALA:O	16:O:108:MET:HG3	2.16	0.44
16:O:158:ILE:O	16:O:176:LYS:HG2	2.17	0.44
1:4:55:ALA:HB1	6:N:137:LEU:HG	2.00	0.44
4:K:201:LEU:O	4:K:205:ALA:N	2.46	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:N:39:ASN:OD1	6:N:39:ASN:N	2.49	0.44
12:L:94:LYS:HE3	12:L:94:LYS:HB3	1.78	0.44
13:B:176:THR:O	13:B:180:ILE:HG12	2.18	0.44
13:B:194:ASP:OD1	13:B:194:ASP:N	2.51	0.44
13:C:156:LEU:HD13	13:C:367:VAL:HG11	2.00	0.44
14:F:35:ILE:HA	14:F:52:VAL:HG12	1.98	0.44
14:F:295:MET:HE3	14:F:295:MET:HB3	1.90	0.44
14:D:139:GLY:HA2	14:D:435:VAL:O	2.17	0.44
16:O:120:VAL:HA	16:O:163:ILE:O	2.18	0.44
13:B:334:VAL:HG21	13:B:351:PHE:CZ	2.53	0.44
6:N:137:LEU:O	6:N:141:LEU:HD13	2.17	0.44
13:A:25:LEU:H	13:A:25:LEU:HD23	1.82	0.44
13:A:307:GLU:HG3	14:E:226:ASN:HB3	2.00	0.44
13:A:382:ALA:N	13:A:487:GLY:O	2.44	0.44
13:B:151:LYS:HD3	13:B:436:MET:HE1	1.99	0.44
13:C:151:LYS:NZ	13:C:465:GLU:OE2	2.47	0.44
14:E:339:SER:HB3	14:E:342:ILE:HG12	1.99	0.44
14:D:119:ILE:HG22	14:D:120:HIS:N	2.32	0.44
15:G:139:GLY:O	15:G:143:VAL:HG23	2.17	0.44
1:4:33:ILE:HG23	1:5:46:LEU:HD22	1.98	0.44
1:5:36:TYR:CE2	1:6:42:LEU:HD13	2.52	0.44
1:8:59:ALA:O	1:8:62:LEU:HB3	2.16	0.44
2:H:95:GLU:OE1	3:I:16:SER:HB2	2.17	0.44
2:H:105:LEU:HD21	2:H:146:GLU:HB2	2.00	0.44
13:A:430:GLN:NE2	13:A:434:SER:O	2.46	0.44
13:C:347:ASP:OD2	14:D:194:ARG:NE	2.47	0.44
14:E:324:ALA:HB3	14:E:325:PRO:HD3	2.00	0.44
1:5:15:THR:HB	1:6:14:ALA:HA	2.00	0.44
1:5:65:LEU:HD11	1:6:63:PHE:HB3	2.00	0.44
13:A:389:THR:HG23	13:A:449:VAL:HG21	2.00	0.44
13:C:247:PRO:HG2	13:C:274:GLN:CD	2.42	0.44
14:E:15:ARG:HH22	16:O:6:ARG:HD3	1.82	0.44
2:H:54:THR:HB	15:G:194:ASP:HB3	2.00	0.44
4:K:90:LEU:HD13	6:N:38:ILE:HD11	2.00	0.44
5:M:33:TRP:O	5:M:34:ASN:C	2.61	0.44
13:A:474:SER:HB3	13:A:475:GLN:HE22	1.83	0.44
14:E:87:ILE:HD13	14:E:238:THR:HG23	1.99	0.44
2:H:143:LYS:HB3	2:H:143:LYS:HE2	1.87	0.43
14:E:85:ILE:O	14:E:119:ILE:HG23	2.17	0.43
14:E:246:PHE:O	14:E:250:GLU:HB2	2.19	0.43
15:G:262:LEU:HD23	15:G:262:LEU:HA	1.85	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:N:30:LEU:HD23	6:N:30:LEU:HA	1.77	0.43
15:G:63:LYS:HB3	15:G:63:LYS:HE3	1.84	0.43
16:O:166:ILE:HG23	16:O:169:LYS:HB2	2.00	0.43
1:8:27:THR:HA	1:1:27:THR:HG21	2.00	0.43
2:H:116:GLN:O	2:H:120:VAL:HG22	2.18	0.43
4:K:89:LYS:O	4:K:93:LEU:HD23	2.19	0.43
5:M:19:ILE:HB	5:M:24:LYS:NZ	2.33	0.43
13:A:270:ASP:OD1	13:A:273:LYS:HD2	2.19	0.43
8:Q:3:GLN:N	8:Q:3:GLN:OE1	2.51	0.43
12:L:61:VAL:HG13	12:L:64:SER:H	1.83	0.43
13:A:64:LEU:HD23	13:A:64:LEU:HA	1.84	0.43
13:A:180:ILE:HD11	13:A:216:LEU:HD11	2.00	0.43
14:E:299:ILE:HD13	14:E:309:SER:HB2	2.01	0.43
14:D:172:LEU:HD23	14:D:172:LEU:HA	1.92	0.43
14:D:422:GLN:O	14:D:425:GLU:HB2	2.18	0.43
15:G:136:PRO:HG2	15:G:218:LYS:HD2	1.99	0.43
1:4:9:ILE:HG13	1:5:6:ALA:HB1	1.99	0.43
5:M:28:SER:O	5:M:31:LYS:HB3	2.18	0.43
10:S:33:ALA:HB1	10:S:37:LEU:HD23	2.00	0.43
13:C:307:GLU:HG3	14:D:226:ASN:HB3	2.00	0.43
1:4:32:LEU:HA	1:3:33:ILE:HG21	2.01	0.43
5:M:19:ILE:HG22	5:M:24:LYS:HD2	2.00	0.43
5:M:54:ALA:HA	5:M:57:LYS:HB2	2.00	0.43
12:L:79:LYS:HB3	12:L:79:LYS:HE3	1.78	0.43
13:A:190:ASN:HA	13:A:198:LYS:HG2	2.00	0.43
13:C:180:ILE:HD12	13:C:220:LEU:HD21	1.99	0.43
14:D:438:LYS:HA	14:D:441:ILE:HG22	2.00	0.43
15:G:50:ALA:HB1	15:G:54:LYS:NZ	2.34	0.43
2:H:40:THR:HG23	2:H:56:GLN:HB3	2.00	0.43
10:S:94:LYS:HD3	10:S:100:TYR:CE1	2.53	0.43
14:E:259:ASP:HA	14:E:260:ASN:HA	1.71	0.43
16:O:163:ILE:HD12	16:O:172:ASP:HA	2.00	0.43
1:4:12:GLY:HA2	1:5:14:ALA:HB2	2.00	0.43
4:K:36:VAL:HA	4:K:39:THR:HG22	2.01	0.43
5:M:155:GLN:NE2	8:Q:33:TYR:O	2.49	0.43
6:N:217:LEU:HD23	6:N:217:LEU:HA	1.67	0.43
14:D:245:TYR:CE1	14:D:249:GLN:HG2	2.54	0.43
15:G:74:ASP:OD1	15:G:110:ASP:N	2.33	0.43
4:K:49:GLU:HG2	9:R:92:TYR:OH	2.19	0.43
5:M:130:MET:HE2	5:M:134:ASP:HB3	2.01	0.43
13:A:271:LEU:HD23	13:A:271:LEU:HA	1.77	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:K:141:ARG:CZ	5:M:46:GLU:HG2	2.48	0.43
6:N:123:ASN:HB2	6:N:127:HIS:CE1	2.54	0.43
11:T:29:ARG:HG2	11:T:33:LEU:HD23	2.01	0.43
12:L:56:THR:OG1	12:L:59:GLY:O	2.36	0.43
14:E:411:ARG:HD2	14:E:415:ARG:HH12	1.84	0.43
2:H:104:ASP:OD1	2:H:104:ASP:N	2.50	0.42
5:M:22:ASN:OD1	5:M:23:GLN:HG2	2.18	0.42
8:Q:43:LYS:HD2	8:Q:43:LYS:HA	1.78	0.42
9:R:27:SER:O	9:R:30:LEU:HG	2.19	0.42
13:A:34:ILE:HD12	13:A:39:ALA:HB2	2.01	0.42
14:E:210:ASN:ND2	14:E:213:ASP:HB2	2.34	0.42
1:1:38:ARG:HG3	1:1:39:ASN:N	2.35	0.42
1:2:37:ALA:HB2	1:3:46:LEU:HD21	2.01	0.42
7:P:3:GLN:O	7:P:5:ILE:HD12	2.19	0.42
14:E:290:THR:O	14:E:294:THR:HG23	2.19	0.42
14:D:52:VAL:HG22	14:D:61:VAL:HG11	2.01	0.42
14:D:171:GLU:HG2	14:D:421:PHE:CD1	2.54	0.42
16:O:148:LEU:O	16:O:149:LYS:HE2	2.19	0.42
2:H:69:ASP:OD1	2:H:70:GLY:N	2.52	0.42
2:H:79:SER:OG	2:H:80:GLY:N	2.50	0.42
6:N:154:MET:O	6:N:158:VAL:HG23	2.20	0.42
13:A:283:LEU:HD13	14:D:279:PRO:HD2	2.02	0.42
2:H:60:PRO:HA	2:H:77:VAL:O	2.20	0.42
3:I:39:SER:O	3:I:39:SER:OG	2.28	0.42
3:I:46:VAL:HG21	15:G:129:LYS:HD3	2.00	0.42
7:P:6:ILE:O	7:P:11:ILE:HG13	2.18	0.42
13:A:131:LEU:HD23	13:A:131:LEU:HA	1.80	0.42
13:A:439:GLU:HG3	13:A:480:LEU:HB3	2.01	0.42
13:B:439:GLU:HG2	13:B:440:GLU:N	2.33	0.42
14:D:90:GLY:HA2	14:D:245:TYR:CE2	2.54	0.42
14:D:152:GLY:HA2	14:D:307:ILE:O	2.19	0.42
16:O:39:LEU:HA	16:O:42:VAL:HG12	2.01	0.42
3:I:37:LYS:HE3	3:I:37:LYS:HB3	1.84	0.42
6:N:186:LEU:N	6:N:187:PRO:HD2	2.33	0.42
13:A:48:GLN:HB2	13:A:51:GLU:HB2	2.01	0.42
14:F:119:ILE:HA	14:F:241:THR:OG1	2.20	0.42
14:F:190:GLY:O	14:F:263:ARG:HG3	2.19	0.42
5:M:109:ALA:O	5:M:112:VAL:HG12	2.19	0.42
6:N:14:ILE:HG12	6:N:15:LEU:HD13	2.01	0.42
6:N:112:THR:HG22	6:N:225:ASN:ND2	2.33	0.42
9:R:26:PRO:HD3	10:S:65:LYS:HZ1	1.84	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
13:B:201:CYS:O	13:B:229:THR:HA	2.20	0.42
13:C:30:ARG:HH11	16:O:58:PRO:HG2	1.85	0.42
14:D:104:PRO:HG3	14:D:110:PRO:HA	2.02	0.42
4:K:60:ALA:HA	6:N:89:PRO:HG3	2.02	0.42
5:M:53:TRP:CZ2	5:M:70:GLU:HB3	2.54	0.42
13:B:450:ARG:HD3	13:B:450:ARG:HA	1.85	0.42
13:C:136:ILE:HG12	14:D:193:THR:HG23	2.01	0.42
16:O:41:ARG:HA	16:O:41:ARG:HD2	1.79	0.42
1:6:62:LEU:HD11	6:N:162:ALA:O	2.19	0.42
4:K:32:THR:O	4:K:36:VAL:HG23	2.20	0.42
4:K:141:ARG:NH2	5:M:46:GLU:HG2	2.35	0.42
10:S:38:VAL:H	11:T:12:LYS:NZ	2.16	0.42
14:F:85:ILE:O	14:F:119:ILE:HG23	2.20	0.42
14:D:245:TYR:O	14:D:249:GLN:HB3	2.20	0.42
16:O:165:ARG:HB2	16:O:170:TYR:HD2	1.85	0.42
13:A:84:GLU:HG3	14:D:32:LEU:HD23	2.02	0.42
13:C:237:SER:HB3	14:F:297:GLU:HG3	2.02	0.42
14:E:145:LEU:HD21	14:E:377:VAL:HG21	2.02	0.42
14:D:160:GLY:HA2	19:D:501:ADP:H1'	2.02	0.42
8:Q:40:LYS:HG3	8:Q:42:MET:H	1.85	0.42
13:A:381:ARG:HH11	13:A:384:LYS:HD2	1.85	0.42
13:A:463:LYS:HB3	13:A:463:LYS:HE2	1.69	0.42
14:D:295:MET:HE3	14:D:295:MET:HB3	1.98	0.42
14:D:441:ILE:HA	14:D:441:ILE:HD12	1.80	0.42
16:O:178:LYS:HD3	16:O:178:LYS:HA	1.79	0.42
4:K:41:LEU:HD12	4:K:41:LEU:HA	1.86	0.41
5:M:29:SER:O	5:M:30:LEU:C	2.63	0.41
13:A:356:LEU:HD23	13:A:359:LYS:HD3	2.01	0.41
14:F:473:ALA:HA	14:F:476:LEU:HD23	2.02	0.41
15:G:10:LEU:HD12	15:G:246:LEU:HB3	2.02	0.41
16:O:16:GLY:O	16:O:20:THR:HG22	2.20	0.41
16:O:122:SER:HA	16:O:158:ILE:HD11	2.02	0.41
1:5:59:ALA:O	1:5:62:LEU:HB3	2.20	0.41
4:K:51:TYR:HA	9:R:77:TYR:OH	2.20	0.41
4:K:125:PHE:HA	4:K:128:GLN:CD	2.45	0.41
6:N:113:VAL:HA	6:N:225:ASN:ND2	2.36	0.41
6:N:120:LYS:HB2	6:N:120:LYS:HE3	1.90	0.41
10:S:49:ILE:O	10:S:53:LYS:NZ	2.51	0.41
13:A:418:LEU:O	13:A:422:VAL:HG23	2.20	0.41
13:A:419:SER:O	13:A:423:ARG:HD3	2.20	0.41
14:E:92:GLU:HB2	14:E:113:THR:HG22	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
14:D:169:ILE:HG23	14:D:257:PHE:CE1	2.55	0.41
15:G:218:LYS:HD2	15:G:218:LYS:HA	1.86	0.41
1:7:72:LEU:HD12	1:7:72:LEU:HA	1.93	0.41
6:N:72:LEU:HD23	6:N:72:LEU:HA	1.74	0.41
8:Q:50:PRO:N	8:Q:51:TRP:HE3	2.18	0.41
9:R:21:LYS:HG3	10:S:70:LYS:HD3	2.02	0.41
13:A:185:ASN:O	13:A:188:ARG:HG2	2.21	0.41
13:A:255:GLU:HG3	13:A:258:ARG:CZ	2.50	0.41
13:B:462:THR:O	13:B:465:GLU:HB3	2.20	0.41
13:C:62:MET:HB2	13:C:76:PHE:CE2	2.55	0.41
13:C:206:ILE:CD1	13:C:247:PRO:HG3	2.50	0.41
14:D:415:ARG:HA	14:D:415:ARG:HD3	1.81	0.41
6:N:73:VAL:O	6:N:76:ILE:HG22	2.20	0.41
12:L:44:VAL:HA	12:L:47:ILE:HG22	2.03	0.41
13:A:69:ASP:OD1	13:A:69:ASP:N	2.40	0.41
13:B:50:GLU:O	13:B:94:ILE:HG23	2.20	0.41
14:D:163:VAL:HG21	14:D:338:LEU:HB3	2.01	0.41
16:O:19:ALA:HB1	16:O:108:MET:SD	2.60	0.41
1:5:42:LEU:HD23	1:5:42:LEU:HA	1.91	0.41
1:1:38:ARG:HB2	1:2:38:ARG:NH2	2.36	0.41
10:S:97:ILE:HG13	10:S:98:ILE:N	2.36	0.41
13:B:156:LEU:HD12	13:B:367:VAL:HG11	2.03	0.41
1:6:74:ALA:HB2	8:Q:1:MET:HB2	2.03	0.41
4:K:103:HIS:HD2	4:K:104:ILE:HD13	1.86	0.41
4:K:154:ARG:HE	12:L:91:PRO:HB3	1.86	0.41
13:A:269:ASP:HA	13:A:270:ASP:HA	1.81	0.41
13:C:170:ASP:O	13:C:175:LYS:NZ	2.54	0.41
14:E:262:PHE:CE2	14:E:316:PRO:HG3	2.56	0.41
14:F:101:ILE:HG13	14:F:103:GLU:HG3	2.03	0.41
14:D:92:GLU:N	14:D:92:GLU:OE1	2.53	0.41
15:G:166:ARG:HE	15:G:166:ARG:HB3	1.63	0.41
4:K:90:LEU:HD12	4:K:90:LEU:HA	1.82	0.41
6:N:60:MET:HE3	6:N:60:MET:HB2	1.89	0.41
7:P:13:MET:HE3	7:P:13:MET:HB3	1.90	0.41
13:B:397:TYR:CD2	13:B:421:GLY:HA3	2.56	0.41
14:D:140:ILE:HB	14:D:143:VAL:HG22	2.03	0.41
15:G:43:ILE:HD13	15:G:43:ILE:HA	1.91	0.41
13:A:44:LEU:HB3	13:A:47:VAL:HB	2.02	0.41
13:A:303:SER:HB2	14:E:225:MET:HG2	2.03	0.41
13:B:469:LEU:O	13:B:473:VAL:HG23	2.21	0.41
13:C:64:LEU:HA	13:C:64:LEU:HD23	1.76	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
14:F:32:LEU:HD12	14:F:32:LEU:HA	1.92	0.41
14:F:348:TYR:HA	14:F:349:PRO:C	2.44	0.41
4:K:157:TYR:O	4:K:160:SER:OG	2.30	0.41
4:K:168:LYS:HD3	4:K:168:LYS:HA	1.84	0.41
8:Q:28:MET:O	8:Q:31:THR:N	2.48	0.41
9:R:20:VAL:HG12	9:R:21:LYS:H	1.85	0.41
13:A:306:LEU:HD23	13:A:306:LEU:HA	1.87	0.41
13:A:468:PHE:O	13:A:472:VAL:HG23	2.20	0.41
13:B:219:ARG:HE	13:B:219:ARG:HB2	1.74	0.41
13:B:439:GLU:HA	13:B:442:VAL:HG12	2.02	0.41
13:C:123:SER:O	13:C:123:SER:OG	2.36	0.41
14:E:402:GLU:O	14:E:406:THR:HG23	2.21	0.41
14:D:342:ILE:H	14:D:342:ILE:HG12	1.65	0.41
16:O:135:LYS:HE3	16:O:139:LYS:HD3	2.03	0.41
16:O:163:ILE:HD11	16:O:174:SER:HB2	2.02	0.41
4:K:134:MET:HE3	4:K:134:MET:HB2	1.85	0.41
4:K:172:HIS:CE1	12:L:64:SER:HA	2.56	0.41
6:N:21:VAL:HG22	8:Q:15:PRO:HG3	2.03	0.41
9:R:24:GLU:O	10:S:65:LYS:NZ	2.44	0.41
13:B:362:ARG:HB3	17:B:601:ATP:N6	2.36	0.41
13:C:34:ILE:HD13	13:C:39:ALA:HB2	2.02	0.41
16:O:76:GLU:O	16:O:76:GLU:HG2	2.20	0.41
1:5:47:PHE:HD1	1:6:49:TYR:OH	2.04	0.40
16:O:22:LEU:HD23	16:O:39:LEU:HD21	2.03	0.40
16:O:67:LYS:HA	16:O:67:LYS:HD3	1.89	0.40
1:8:58:GLU:O	1:8:59:ALA:C	2.65	0.40
4:K:39:THR:HA	4:K:42:ILE:HG22	2.02	0.40
5:M:13:VAL:O	5:M:17:GLU:HG2	2.20	0.40
12:L:66:GLU:HG3	12:L:67:TYR:H	1.85	0.40
13:A:460:LYS:HD3	13:A:508:PHE:CE1	2.55	0.40
13:B:347:ASP:OD1	14:F:194:ARG:NE	2.45	0.40
14:F:24:VAL:O	14:F:63:THR:OG1	2.32	0.40
1:5:5:ALA:CB	1:6:6:ALA:HB2	2.48	0.40
2:H:78:SER:HB3	3:I:23:VAL:CG2	2.52	0.40
2:H:124:ASP:HB3	2:H:127:THR:HG22	2.03	0.40
12:L:72:GLU:HG3	12:L:73:ARG:N	2.36	0.40
12:L:83:GLY:O	12:L:84:ASN:ND2	2.55	0.40
13:B:32:LEU:N	13:B:40:ARG:O	2.52	0.40
13:B:175:LYS:HD2	13:B:326:VAL:HG13	2.02	0.40
13:B:242:LEU:HD23	13:B:242:LEU:HA	1.86	0.40
13:C:447:ALA:O	13:C:452:TYR:HB2	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
14:E:97:ILE:HG12	14:E:220:LEU:HB2	2.02	0.40
15:G:63:LYS:HG3	15:G:156:ASP:OD1	2.21	0.40
16:O:135:LYS:O	16:O:139:LYS:HB2	2.21	0.40
10:S:91:ILE:HA	10:S:94:LYS:HE2	2.04	0.40
12:L:90:PHE:CG	12:L:91:PRO:HD2	2.57	0.40
14:D:95:GLY:N	14:D:218:VAL:O	2.32	0.40
14:D:157:LEU:HD23	14:D:336:THR:HB	2.03	0.40
15:G:201:LEU:O	15:G:205:GLN:HG2	2.21	0.40
16:O:153:LYS:HD3	16:O:153:LYS:HA	1.83	0.40
1:4:58:GLU:O	1:4:59:ALA:C	2.65	0.40
4:K:101:ILE:HG23	5:M:118:MET:HE2	2.04	0.40
6:N:41:ARG:CZ	9:R:59:LYS:HA	2.51	0.40
13:A:124:LYS:HE2	13:A:124:LYS:HB2	1.96	0.40
13:A:151:LYS:NZ	13:A:465:GLU:OE1	2.52	0.40
13:A:345:ILE:HA	14:E:225:MET:SD	2.62	0.40
17:B:601:ATP:H8	17:B:601:ATP:H5'1	1.86	0.40
14:E:282:VAL:HG21	14:E:323:PRO:CD	2.51	0.40
14:F:166:THR:N	19:F:501:ADP:O1A	2.55	0.40
14:D:171:GLU:OE2	14:D:175:ASN:ND2	2.54	0.40
14:D:371:TYR:OH	14:D:375:ARG:NH2	2.52	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	1	73/75 (97%)	67 (92%)	6 (8%)	0	100	100
1	2	73/75 (97%)	66 (90%)	7 (10%)	0	100	100
1	3	73/75 (97%)	67 (92%)	6 (8%)	0	100	100
1	4	73/75 (97%)	67 (92%)	6 (8%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	5	73/75 (97%)	66 (90%)	7 (10%)	0	100	100
1	6	73/75 (97%)	66 (90%)	7 (10%)	0	100	100
1	7	73/75 (97%)	67 (92%)	6 (8%)	0	100	100
1	8	73/75 (97%)	67 (92%)	6 (8%)	0	100	100
2	H	130/146 (89%)	121 (93%)	9 (7%)	0	100	100
3	I	43/51 (84%)	41 (95%)	2 (5%)	0	100	100
4	K	194/214 (91%)	186 (96%)	8 (4%)	0	100	100
5	M	154/160 (96%)	146 (95%)	8 (5%)	0	100	100
6	N	221/226 (98%)	204 (92%)	17 (8%)	0	100	100
7	P	39/58 (67%)	36 (92%)	3 (8%)	0	100	100
8	Q	49/68 (72%)	40 (82%)	9 (18%)	0	100	100
9	R	72/93 (77%)	68 (94%)	4 (6%)	0	100	100
10	S	75/102 (74%)	74 (99%)	1 (1%)	0	100	100
11	T	42/69 (61%)	42 (100%)	0	0	100	100
12	L	65/108 (60%)	60 (92%)	5 (8%)	0	100	100
13	A	484/510 (95%)	466 (96%)	18 (4%)	0	100	100
13	B	473/510 (93%)	449 (95%)	24 (5%)	0	100	100
13	C	466/510 (91%)	442 (95%)	24 (5%)	0	100	100
14	D	424/482 (88%)	382 (90%)	42 (10%)	0	100	100
14	E	452/482 (94%)	430 (95%)	21 (5%)	1 (0%)	44	76
14	F	464/482 (96%)	434 (94%)	30 (6%)	0	100	100
15	G	269/273 (98%)	259 (96%)	10 (4%)	0	100	100
16	O	185/190 (97%)	177 (96%)	8 (4%)	0	100	100
All	All	4885/5334 (92%)	4590 (94%)	294 (6%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
14	E	280	SER

5.3.2 Protein sidechains (i)

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	1	51/51 (100%)	50 (98%)	1 (2%)	50 77
1	2	51/51 (100%)	51 (100%)	0	100 100
1	3	51/51 (100%)	51 (100%)	0	100 100
1	4	51/51 (100%)	51 (100%)	0	100 100
1	5	51/51 (100%)	51 (100%)	0	100 100
1	6	51/51 (100%)	51 (100%)	0	100 100
1	7	51/51 (100%)	51 (100%)	0	100 100
1	8	51/51 (100%)	51 (100%)	0	100 100
2	H	103/108 (95%)	103 (100%)	0	100 100
3	I	36/42 (86%)	36 (100%)	0	100 100
4	K	160/187 (86%)	160 (100%)	0	100 100
5	M	135/139 (97%)	135 (100%)	0	100 100
6	N	196/199 (98%)	196 (100%)	0	100 100
7	P	36/48 (75%)	36 (100%)	0	100 100
8	Q	51/68 (75%)	51 (100%)	0	100 100
9	R	65/82 (79%)	65 (100%)	0	100 100
10	S	64/85 (75%)	64 (100%)	0	100 100
11	T	34/58 (59%)	34 (100%)	0	100 100
12	L	60/98 (61%)	60 (100%)	0	100 100
13	A	393/412 (95%)	393 (100%)	0	100 100
13	B	387/412 (94%)	387 (100%)	0	100 100
13	C	382/412 (93%)	381 (100%)	1 (0%)	91 96
14	D	348/388 (90%)	347 (100%)	1 (0%)	91 96
14	E	369/388 (95%)	369 (100%)	0	100 100
14	F	376/388 (97%)	376 (100%)	0	100 100
15	G	228/229 (100%)	228 (100%)	0	100 100
16	O	162/165 (98%)	162 (100%)	0	100 100
All	All	3993/4316 (92%)	3990 (100%)	3 (0%)	92 97

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	1	38	ARG
13	C	373	ARG
14	D	158	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (66) such sidechains are listed below:

Mol	Chain	Res	Type
1	2	45	GLN
2	H	85	ASN
2	H	132	GLN
2	H	138	ASN
3	I	17	GLN
3	I	34	ASN
4	K	103	HIS
4	K	105	GLN
4	K	114	GLN
4	K	122	HIS
4	K	130	ASN
4	K	153	ASN
4	K	163	ASN
5	M	23	GLN
5	M	34	ASN
6	N	46	GLN
6	N	47	GLN
6	N	61	HIS
6	N	101	ASN
6	N	225	ASN
8	Q	32	ASN
8	Q	46	ASN
8	Q	48	ASN
9	R	44	GLN
9	R	83	HIS
10	S	57	ASN
12	L	84	ASN
12	L	88	ASN
13	A	349	GLN
13	A	385	GLN
13	A	396	GLN
13	A	415	GLN
13	A	471	HIS
13	A	475	GLN
13	B	65	ASN
13	B	471	HIS

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Mol	Chain	Res	Type
13	B	475	GLN
13	B	492	GLN
13	C	172	GLN
13	C	208	GLN
13	C	385	GLN
13	C	471	HIS
14	E	210	ASN
14	E	224	GLN
14	E	311	GLN
14	E	370	HIS
14	E	382	GLN
14	E	388	GLN
14	F	133	GLN
14	F	249	GLN
14	F	252	GLN
14	F	311	GLN
14	F	331	HIS
14	F	364	ASN
14	D	76	GLN
14	D	197	ASN
14	D	249	GLN
14	D	378	GLN
14	D	382	GLN
14	D	419	GLN
15	G	15	ASN
15	G	66	HIS
15	G	88	GLN
15	G	120	HIS
15	G	163	ASN
16	O	84	ASN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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 10 ligands modelled in this entry, 5 are monoatomic - leaving 5 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	$\#\|Z\| > 2$
19	ADP	F	501	-	24,29,29	2.38	8 (33%)	29,45,45	1.59	4 (13%)
17	ATP	A	601	18	26,33,33	2.24	6 (23%)	31,52,52	1.72	6 (19%)
17	ATP	B	601	18	26,33,33	2.22	7 (26%)	31,52,52	1.76	7 (22%)
17	ATP	C	601	18	26,33,33	2.23	6 (23%)	31,52,52	1.59	7 (22%)
19	ADP	D	501	18	24,29,29	2.31	8 (33%)	29,45,45	1.73	6 (20%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
19	ADP	F	501	-	-	7/12/32/32	0/3/3/3
17	ATP	A	601	18	-	0/18/38/38	0/3/3/3
17	ATP	B	601	18	-	2/18/38/38	0/3/3/3
17	ATP	C	601	18	-	9/18/38/38	0/3/3/3
19	ADP	D	501	18	-	3/12/32/32	0/3/3/3

All (35) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
17	A	601	ATP	O4'-C1'	7.57	1.51	1.41
17	C	601	ATP	O4'-C1'	7.27	1.51	1.41
17	B	601	ATP	O4'-C1'	7.24	1.51	1.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
19	D	501	ADP	O4'-C1'	6.50	1.50	1.41
19	F	501	ADP	O4'-C1'	6.27	1.49	1.41
17	C	601	ATP	C2'-C1'	-4.65	1.46	1.53
17	B	601	ATP	C2'-C1'	-4.53	1.46	1.53
19	D	501	ADP	C6-N6	4.49	1.50	1.34
19	F	501	ADP	C6-N6	4.46	1.50	1.34
19	F	501	ADP	C2'-C1'	-4.46	1.47	1.53
17	A	601	ATP	C2'-C1'	-4.37	1.47	1.53
19	D	501	ADP	C2'-C1'	-3.84	1.47	1.53
17	A	601	ATP	C6-N6	3.43	1.46	1.34
17	B	601	ATP	C2'-C3'	-3.39	1.44	1.53
17	C	601	ATP	C6-N6	3.30	1.46	1.34
17	B	601	ATP	C6-N6	3.26	1.45	1.34
17	C	601	ATP	C2'-C3'	-3.26	1.44	1.53
19	D	501	ADP	O4'-C4'	3.21	1.52	1.45
19	F	501	ADP	O4'-C4'	3.20	1.52	1.45
17	A	601	ATP	C2'-C3'	-3.14	1.44	1.53
19	F	501	ADP	C2'-C3'	-2.94	1.45	1.53
19	F	501	ADP	C3'-C4'	-2.88	1.45	1.53
17	C	601	ATP	C4-N3	-2.64	1.32	1.35
17	B	601	ATP	C4-N3	-2.59	1.32	1.35
17	A	601	ATP	C4-N3	-2.57	1.32	1.35
19	D	501	ADP	C3'-C4'	-2.54	1.46	1.53
19	D	501	ADP	C2'-C3'	-2.50	1.46	1.53
17	A	601	ATP	O4'-C4'	2.43	1.50	1.45
17	C	601	ATP	O4'-C4'	2.32	1.50	1.45
19	D	501	ADP	PA-O5'	2.27	1.68	1.59
17	B	601	ATP	O4'-C4'	2.25	1.50	1.45
19	F	501	ADP	C6-C5	-2.15	1.35	1.43
17	B	601	ATP	C3'-C4'	-2.07	1.47	1.53
19	F	501	ADP	PA-O5'	2.03	1.67	1.59
19	D	501	ADP	C6-C5	-2.02	1.35	1.43

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
17	A	601	ATP	C3'-C2'-C1'	4.52	107.78	100.98
19	D	501	ADP	C3'-C2'-C1'	4.38	107.57	100.98
19	D	501	ADP	N3-C2-N1	-4.13	122.22	128.68
17	B	601	ATP	N3-C2-N1	-4.12	122.23	128.68
17	C	601	ATP	N3-C2-N1	-4.12	122.23	128.68
19	F	501	ADP	C3'-C2'-C1'	4.07	107.11	100.98

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
17	B	601	ATP	C3'-C2'-C1'	4.06	107.09	100.98
19	F	501	ADP	N3-C2-N1	-3.95	122.51	128.68
17	A	601	ATP	N3-C2-N1	-3.70	122.89	128.68
19	D	501	ADP	C4-C5-N7	-3.68	105.57	109.40
17	A	601	ATP	C4-C5-N7	-3.66	105.58	109.40
19	F	501	ADP	C4-C5-N7	-3.62	105.62	109.40
17	C	601	ATP	C4-C5-N7	-3.60	105.64	109.40
17	B	601	ATP	C4-C5-N7	-3.56	105.69	109.40
17	B	601	ATP	PA-O3A-PB	-3.41	121.14	132.83
17	C	601	ATP	C3'-C2'-C1'	3.35	106.02	100.98
17	B	601	ATP	PB-O3B-PG	-3.21	121.81	132.83
19	D	501	ADP	PA-O3A-PB	-3.20	121.85	132.83
17	A	601	ATP	PB-O3B-PG	-3.20	121.85	132.83
17	A	601	ATP	PA-O3A-PB	-2.89	122.90	132.83
17	C	601	ATP	PA-O3A-PB	-2.86	123.01	132.83
19	F	501	ADP	PA-O3A-PB	-2.84	123.07	132.83
19	D	501	ADP	C1'-N9-C4	-2.76	121.80	126.64
17	A	601	ATP	C2'-C3'-C4'	2.72	107.92	102.64
17	C	601	ATP	C2'-C3'-C4'	2.38	107.26	102.64
17	B	601	ATP	C2'-C3'-C4'	2.34	107.20	102.64
17	B	601	ATP	C1'-N9-C4	-2.18	122.81	126.64
17	C	601	ATP	PB-O3B-PG	-2.08	125.69	132.83
19	D	501	ADP	C2'-C3'-C4'	2.07	106.66	102.64
17	C	601	ATP	C1'-N9-C4	-2.03	123.07	126.64

There are no chirality outliers.

All (21) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
17	C	601	ATP	C5'-O5'-PA-O1A
17	C	601	ATP	C5'-O5'-PA-O3A
19	F	501	ADP	PA-O3A-PB-O2B
19	F	501	ADP	PA-O3A-PB-O3B
19	F	501	ADP	C5'-O5'-PA-O2A
19	F	501	ADP	O4'-C4'-C5'-O5'
19	F	501	ADP	C3'-C4'-C5'-O5'
19	D	501	ADP	C3'-C4'-C5'-O5'
17	C	601	ATP	O4'-C4'-C5'-O5'
17	C	601	ATP	C3'-C4'-C5'-O5'
17	B	601	ATP	O4'-C4'-C5'-O5'
19	D	501	ADP	O4'-C4'-C5'-O5'
19	F	501	ADP	C5'-O5'-PA-O3A

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Mol	Chain	Res	Type	Atoms
17	C	601	ATP	PG-O3B-PB-O1B
17	C	601	ATP	C5'-O5'-PA-O2A
19	F	501	ADP	C5'-O5'-PA-O1A
17	C	601	ATP	PB-O3B-PG-O1G
17	C	601	ATP	PB-O3B-PG-O2G
17	B	601	ATP	PB-O3A-PA-O2A
17	C	601	ATP	PG-O3B-PB-O2B
19	D	501	ADP	C5'-O5'-PA-O1A

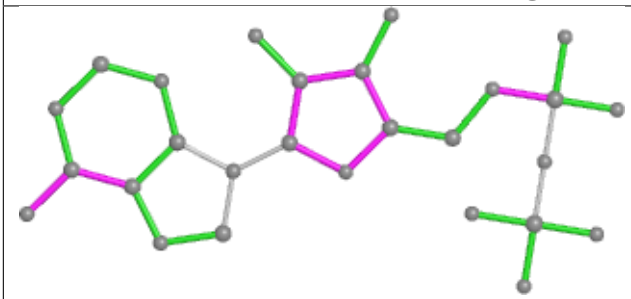
There are no ring outliers.

5 monomers are involved in 9 short contacts:

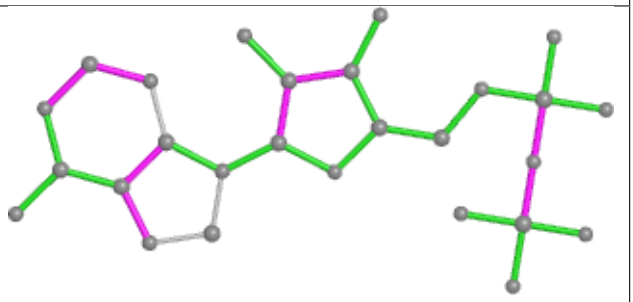
Mol	Chain	Res	Type	Clashes	Symm-Clashes
19	F	501	ADP	1	0
17	A	601	ATP	1	0
17	B	601	ATP	3	0
17	C	601	ATP	1	0
19	D	501	ADP	3	0

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 then 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.

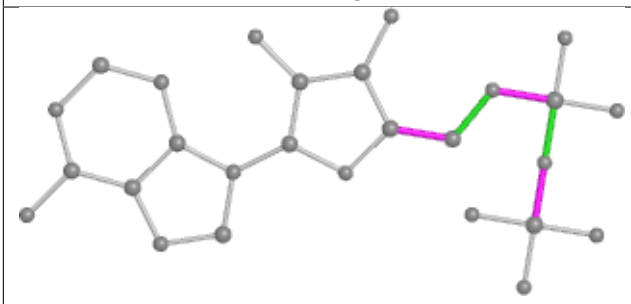
Ligand ADP F 501



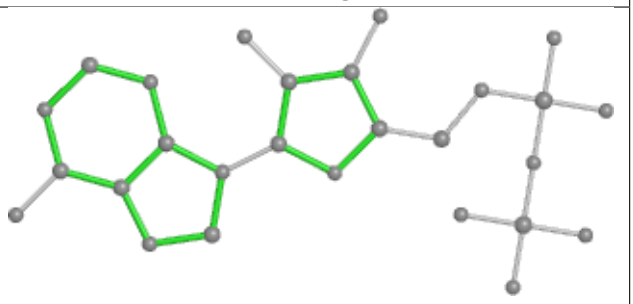
Bond lengths



Bond angles

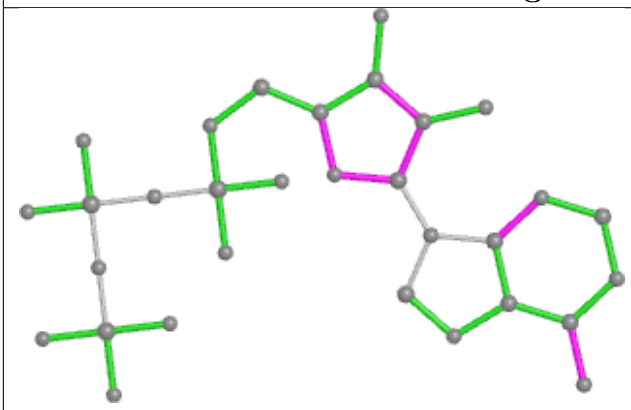


Torsions

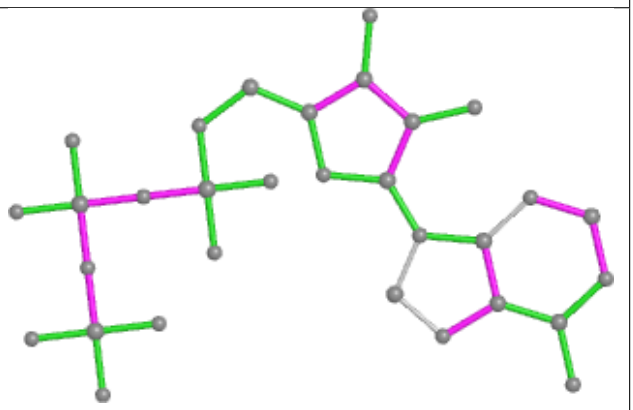


Rings

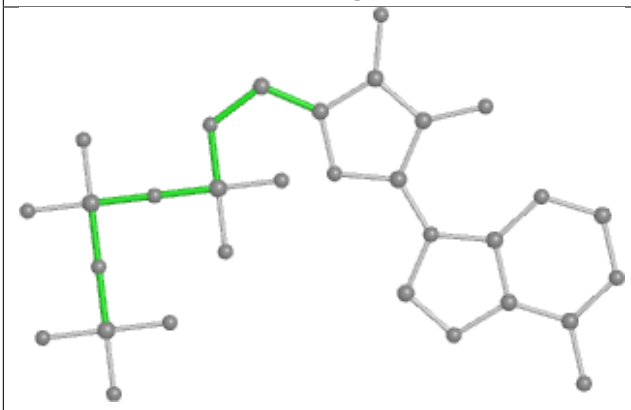
Ligand ATP A 601



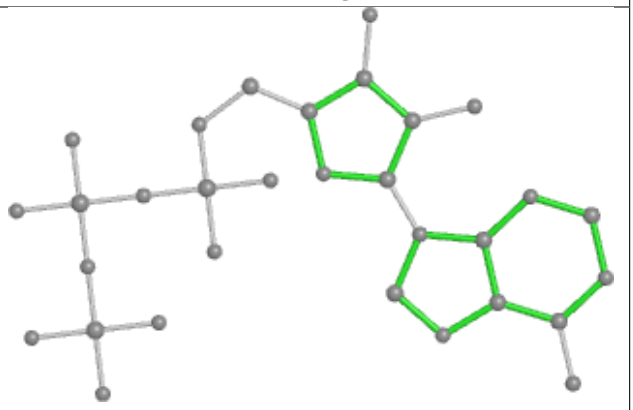
Bond lengths



Bond angles

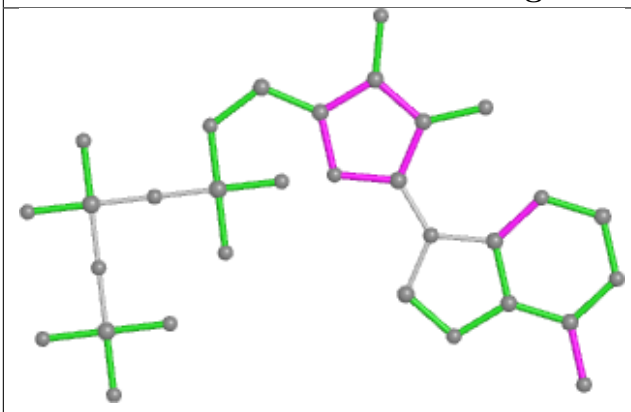


Torsions

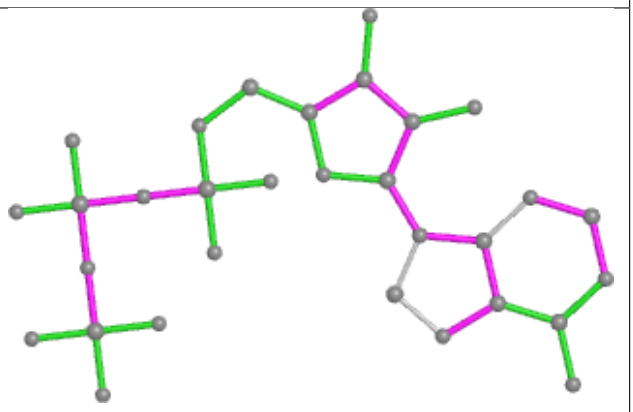


Rings

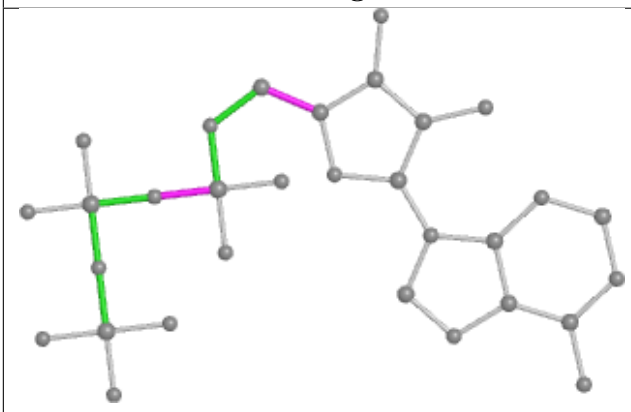
Ligand ATP B 601



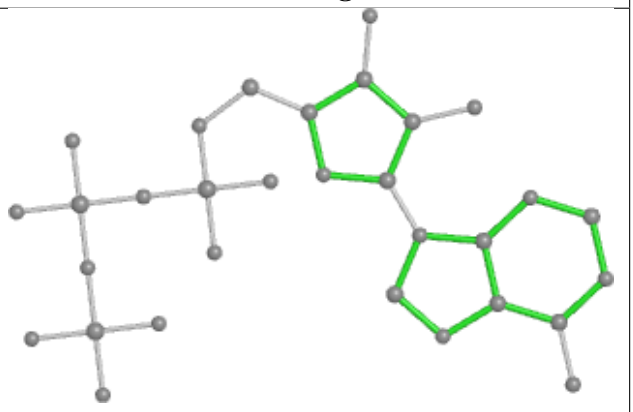
Bond lengths



Bond angles

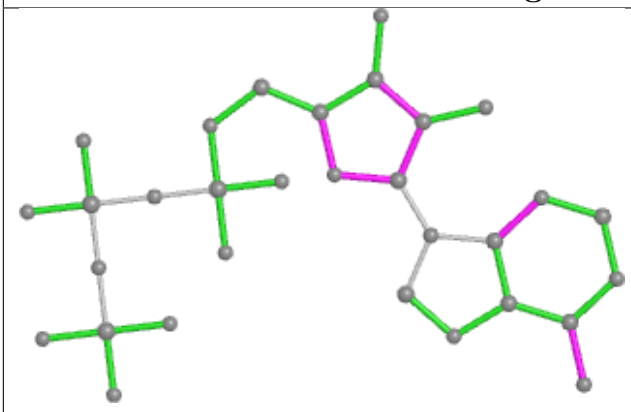


Torsions

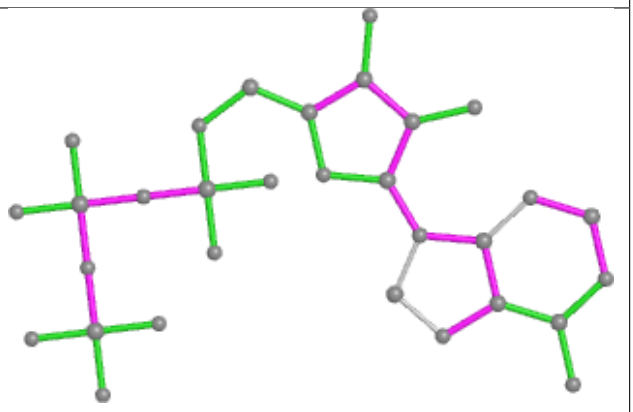


Rings

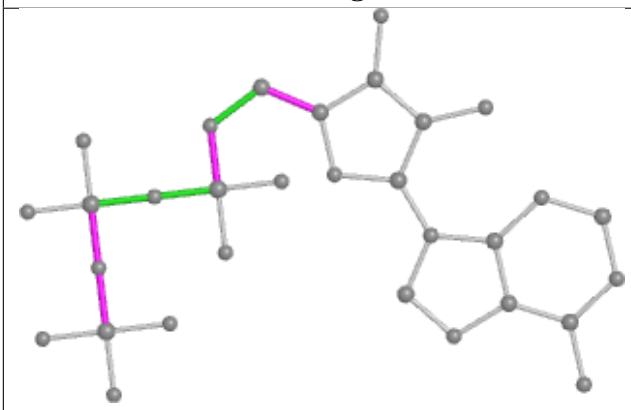
Ligand ATP C 601



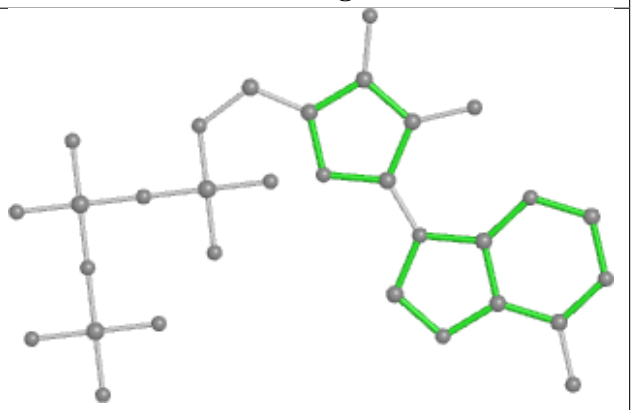
Bond lengths



Bond angles

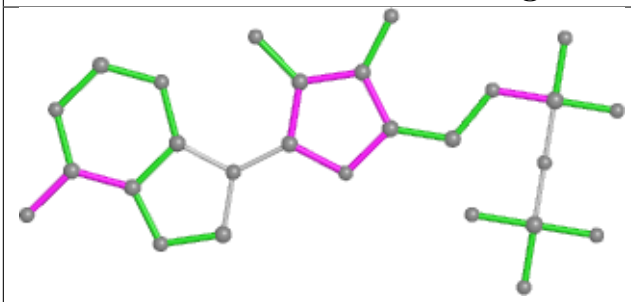


Torsions

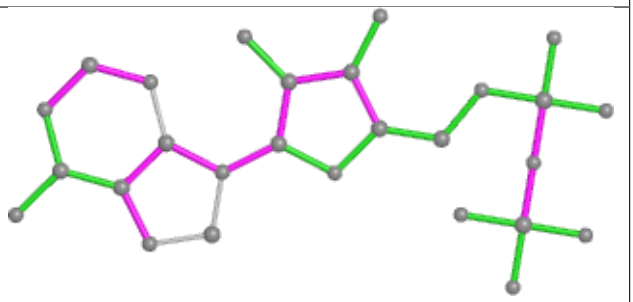


Rings

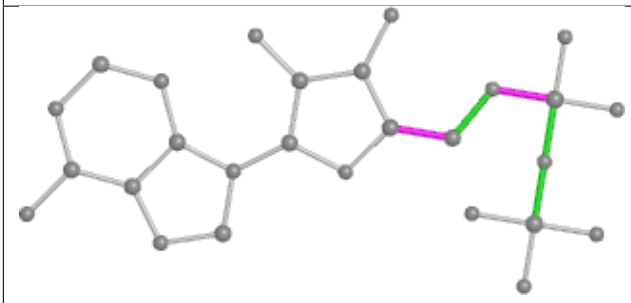
Ligand ADP D 501



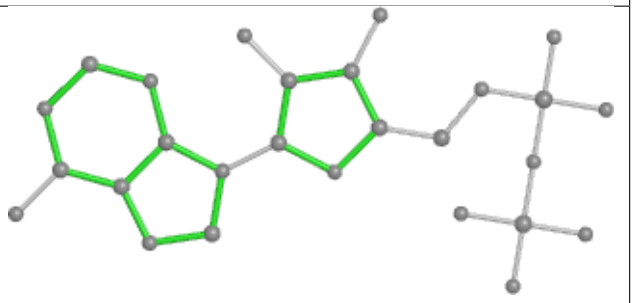
Bond lengths



Bond angles



Torsions



Rings

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.

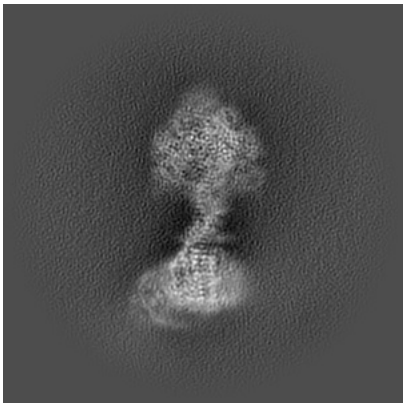
6 Map visualisation

This section contains visualisations of the EMDB entry EMD-34583. These allow visual inspection of the internal detail of the map and identification of artifacts.

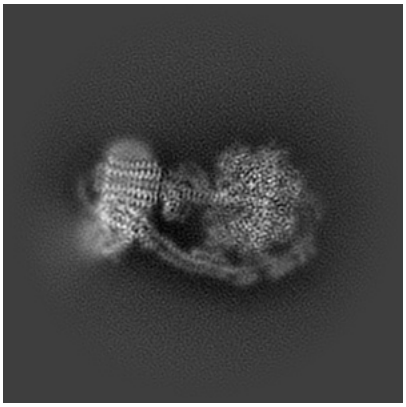
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections

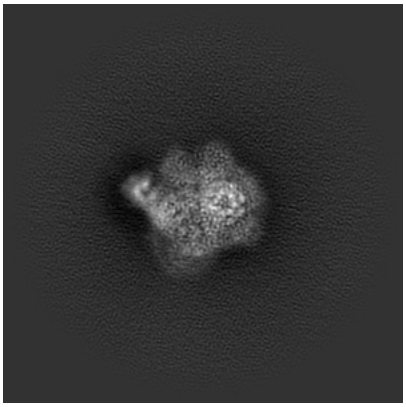
6.1.1 Primary map



X



Y

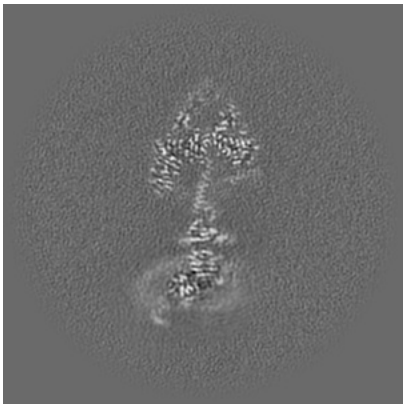


Z

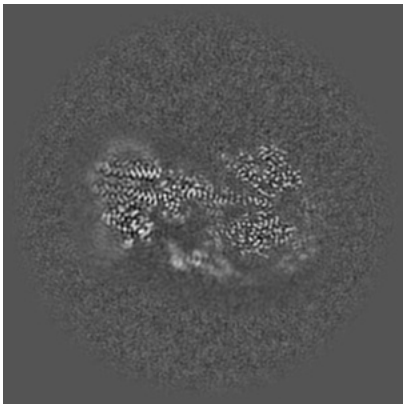
The images above show the map projected in three orthogonal directions.

6.2 Central slices

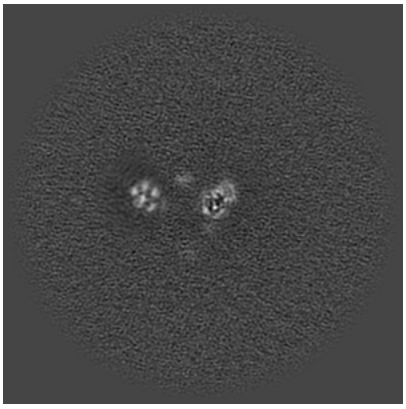
6.2.1 Primary map



X Index: 256



Y Index: 256

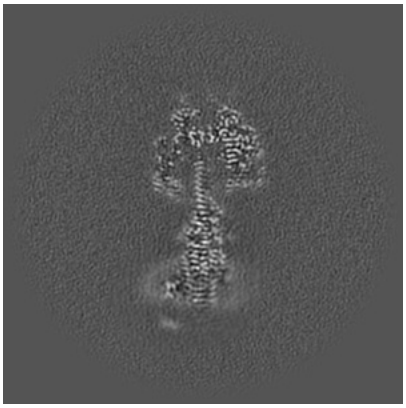


Z Index: 256

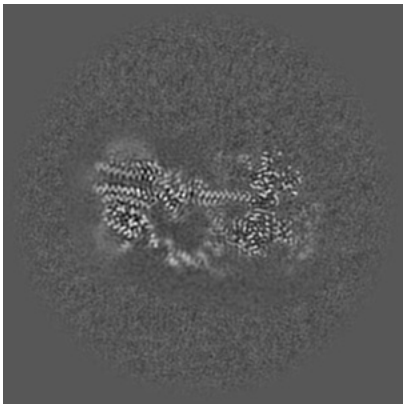
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices (i)

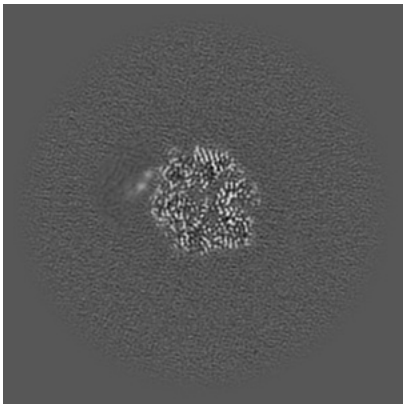
6.3.1 Primary map



X Index: 267



Y Index: 250

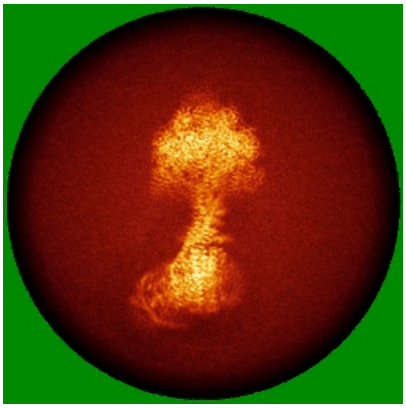


Z Index: 328

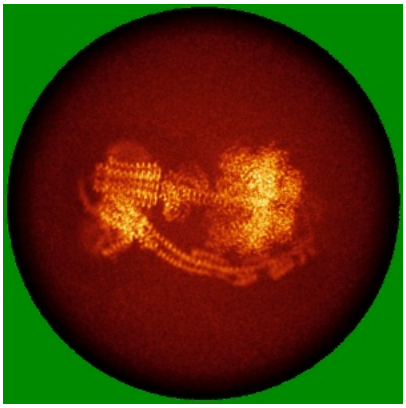
The images above show the largest variance slices of the map in three orthogonal directions.

6.4 Orthogonal standard-deviation projections (False-color) (i)

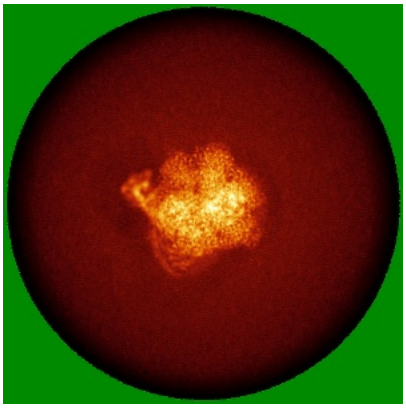
6.4.1 Primary map



X



Y

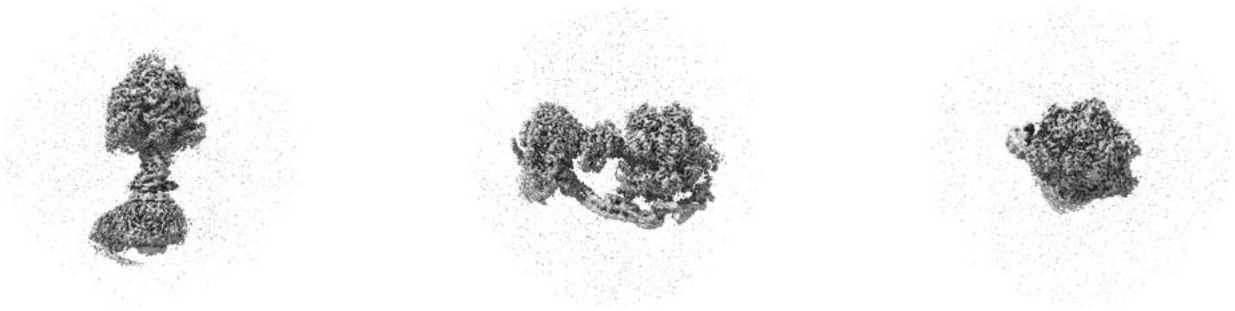


Z

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



X Y Z

The images above show the 3D surface view of the map at the recommended contour level 4.5. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

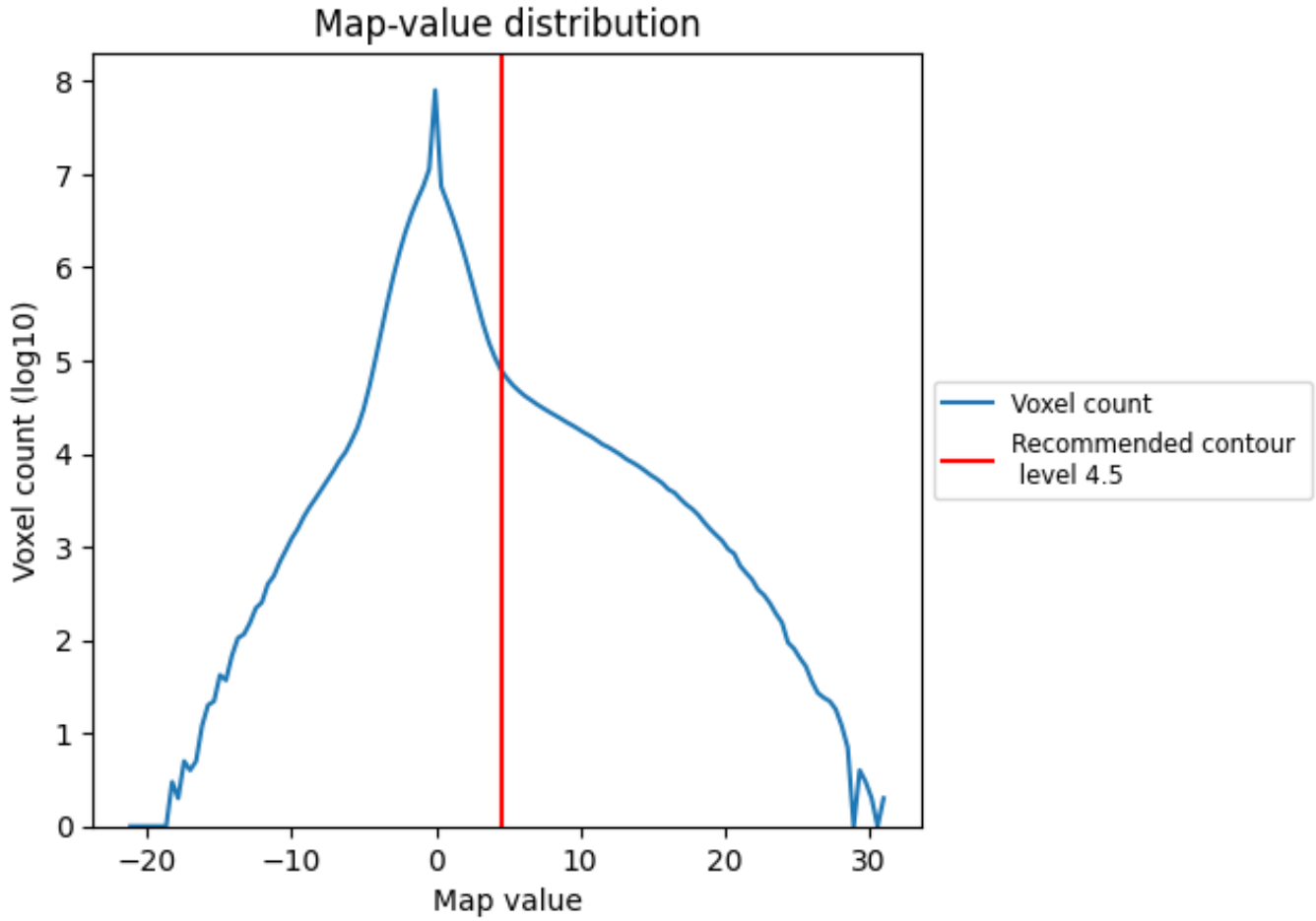
6.6 Mask visualisation (i)

This section was not generated. No masks/segmentation were deposited.

7 Map analysis

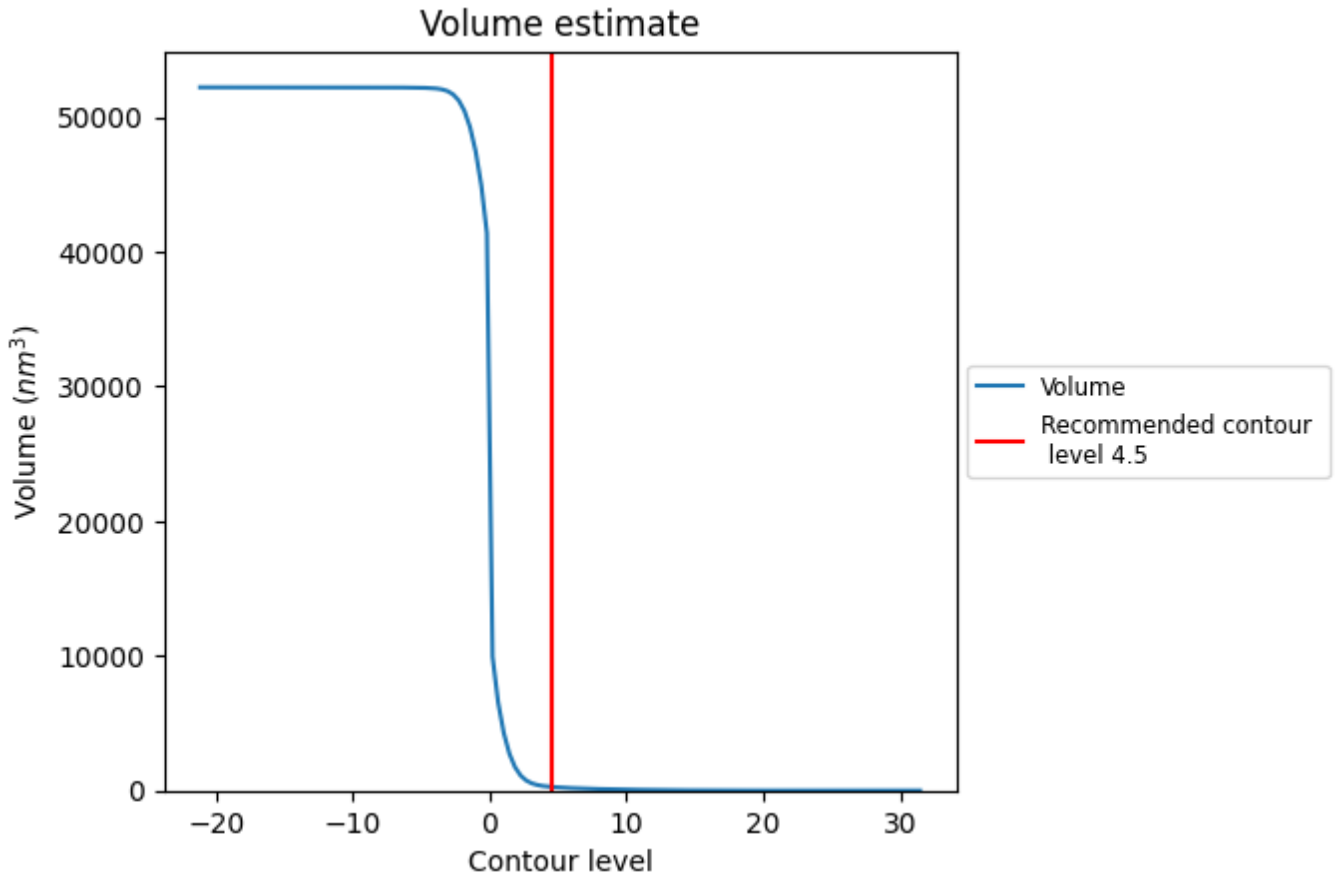
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution



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.

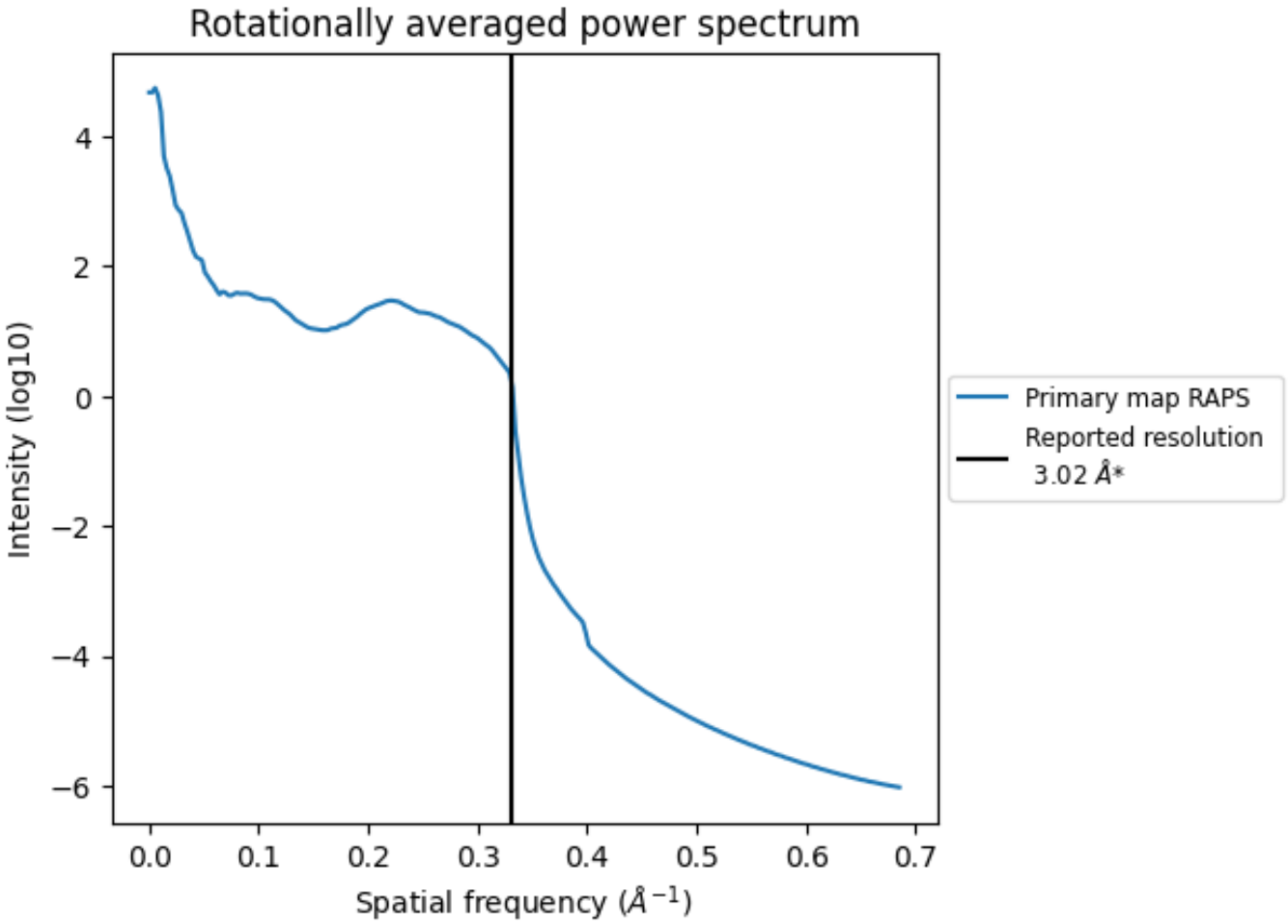
7.2 Volume estimate



The volume at the recommended contour level is 272 nm^3; this corresponds to an approximate mass of 246 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 (i)



*Reported resolution corresponds to spatial frequency of 0.331 Å^{-1}

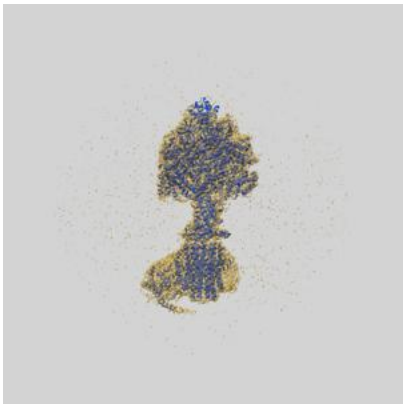
8 Fourier-Shell correlation (i)

This section was not generated. No FSC curve or half-maps provided.

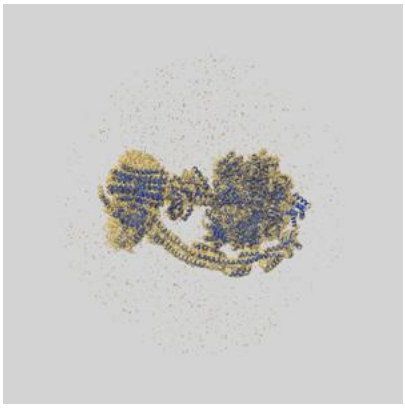
9 Map-model fit

This section contains information regarding the fit between EMDB map EMD-34583 and PDB model 8H9V. Per-residue inclusion information can be found in section 3 on page 9.

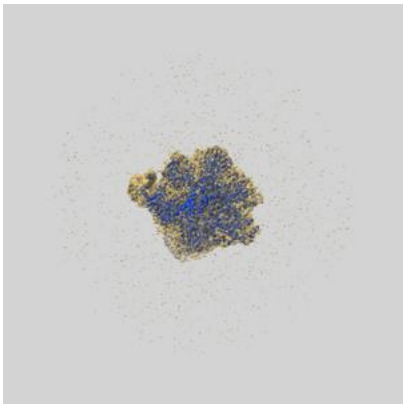
9.1 Map-model overlay



X



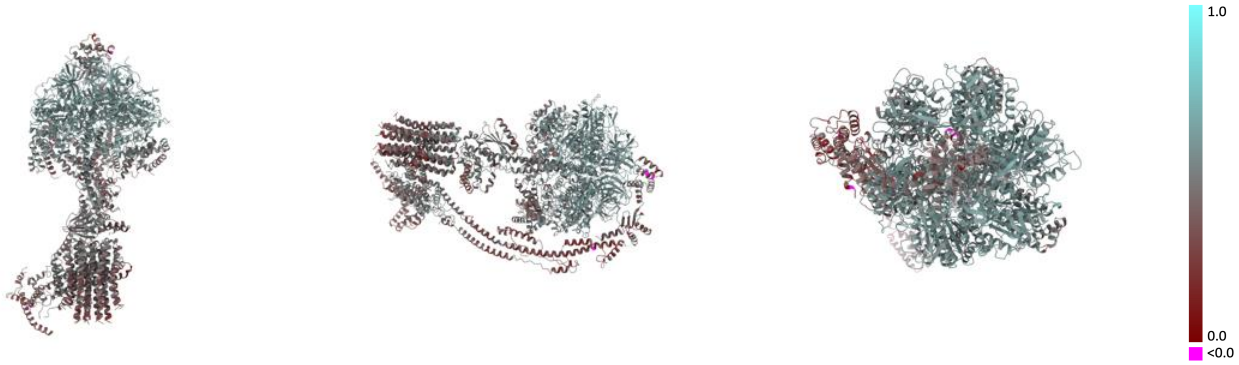
Y



Z

The images above show the 3D surface view of the map at the recommended contour level 4.5 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.

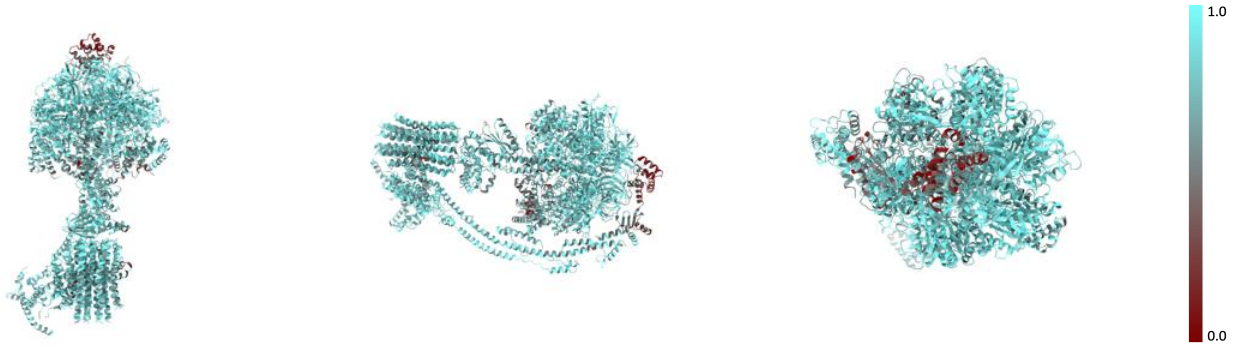
9.2 Q-score mapped to coordinate model (i)



X Y Z

The images above show the model with each residue coloured according 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.

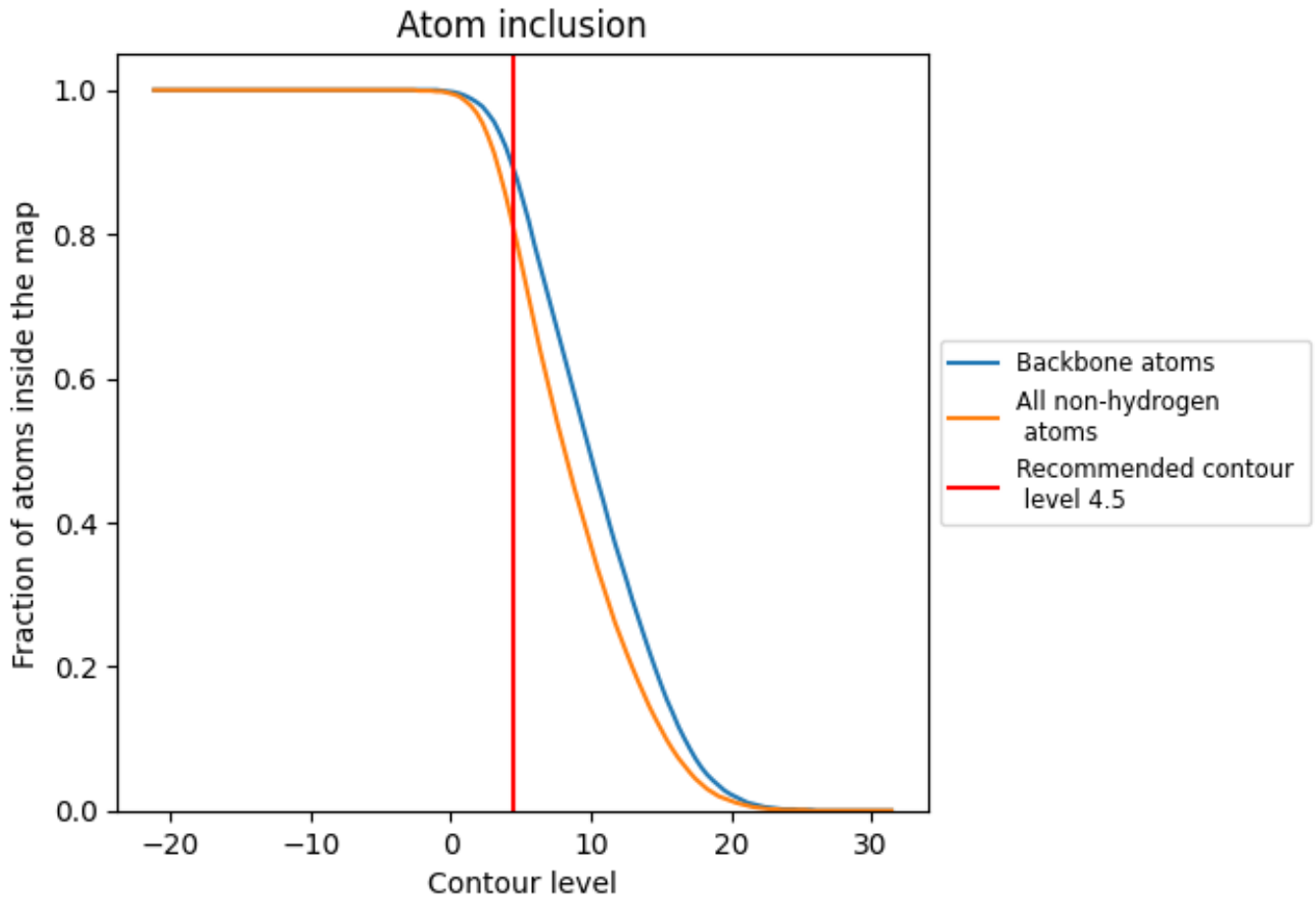
9.3 Atom inclusion mapped to coordinate model (i)



X Y Z

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 (4.5).

9.4 Atom inclusion (i)



At the recommended contour level, 89% of all backbone atoms, 81% 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 (4.5) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.8060	0.4870
1	0.8210	0.4000
2	0.8260	0.4170
3	0.7740	0.3940
4	0.7790	0.3950
5	0.7830	0.3920
6	0.8190	0.4140
7	0.7700	0.3830
8	0.7790	0.3980
A	0.8000	0.5400
B	0.8620	0.5610
C	0.8690	0.5570
D	0.7470	0.5140
E	0.8550	0.5580
F	0.8720	0.5670
G	0.7520	0.4730
H	0.8360	0.4470
I	0.7940	0.4600
K	0.8340	0.3700
L	0.7820	0.2690
M	0.8550	0.3640
N	0.8530	0.4650
O	0.3510	0.3970
P	0.8120	0.3970
Q	0.8400	0.4370
R	0.8880	0.4270
S	0.7240	0.3140
T	0.7260	0.3210