Full wwPDB EM Validation Report ⓘ

Jun 10, 2025 – 03:08 PM JST

PDB ID	:	8H9I / pdb_00008h9i
EMDB ID	:	EMD-34568
Title	:	Human ATP synthase F1 domain, state2
Authors	:	Lai, Y.; Zhang, Y.; Liu, F.; Gao, Y.; Gong, H.; Rao, Z.
Deposited on	:	2022-10-25
Resolution	:	2.77 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:

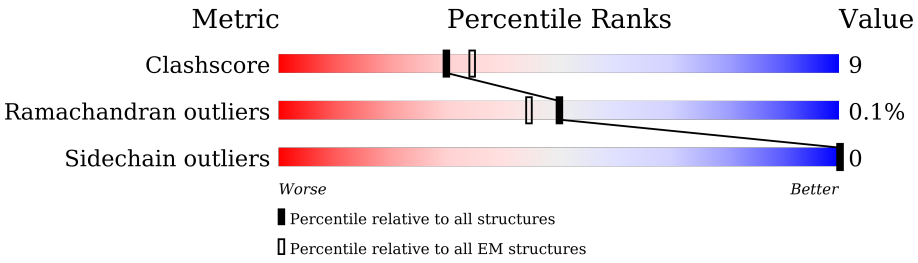
EMDB validation analysis	:	0.0.1.dev118
Mogul	:	1.8.5 (274361), CSD as541be (2020)
MolProbity	:	4-5-2 with Phenix2.0rc1
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
MapQ	:	1.9.13
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.43.1

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 2.77 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	A	510	77% 17% 6%
1	B	510	78% 17% 5%
1	C	510	79% 20%
2	D	482	11% 69% 26% 5%
2	E	482	77% 20%
2	F	482	77% 20%
3	G	273	24% 5% 70%
4	O	190	57% 64% 35%

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 23961 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 subunit alpha, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	C	503	Total	C	N	O	S	0	0
			3830	2410	675	732	13		
1	A	477	Total	C	N	O	S	0	0
			3639	2290	645	692	12		
1	B	487	Total	C	N	O	S	0	0
			3714	2338	656	707	13		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
2	E	466	Total	C	N	O	S	0	0
			3522	2234	597	678	13		
2	F	470	Total	C	N	O	S	0	0
			3562	2257	604	688	13		
2	D	460	Total	C	N	O	S	0	0
			3482	2204	591	674	13		

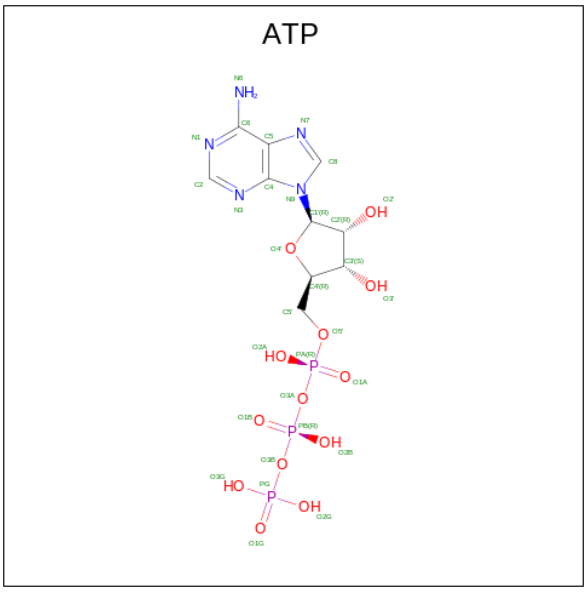
- Molecule 3 is a protein called ATP synthase subunit gamma, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	G	81	Total	C	N	O	S	0	0
			623	387	113	118	5		

- Molecule 4 is a protein called ATP synthase subunit O, mitochondrial.

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

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

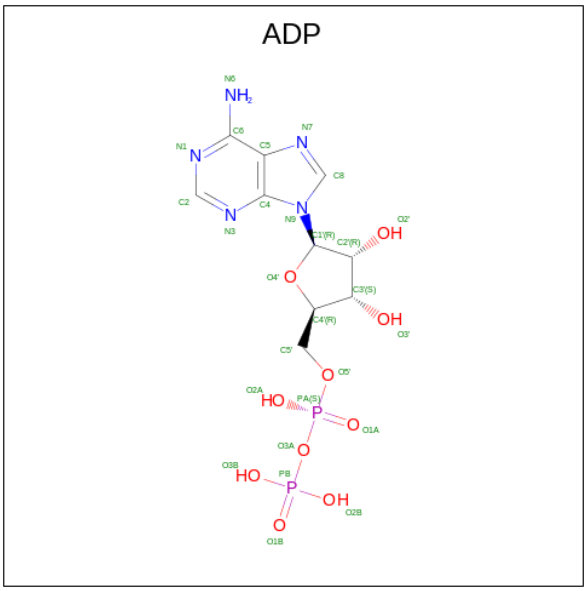


Mol	Chain	Residues	Atoms					AltConf
5	C	1	Total	C	N	O	P	0
			31	10	5	13	3	
5	A	1	Total	C	N	O	P	0
			31	10	5	13	3	
5	B	1	Total	C	N	O	P	0
			31	10	5	13	3	

- Molecule 6 is MAGNESIUM ION (CCD ID: MG) (formula: Mg) (labeled as "Ligand of Interest" by depositor).

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

- Molecule 7 is ADENOSINE-5'-DIPHOSPHATE (CCD ID: ADP) (formula: $C_{10}H_{15}N_5O_{10}P_2$) (labeled as "Ligand of Interest" by depositor).

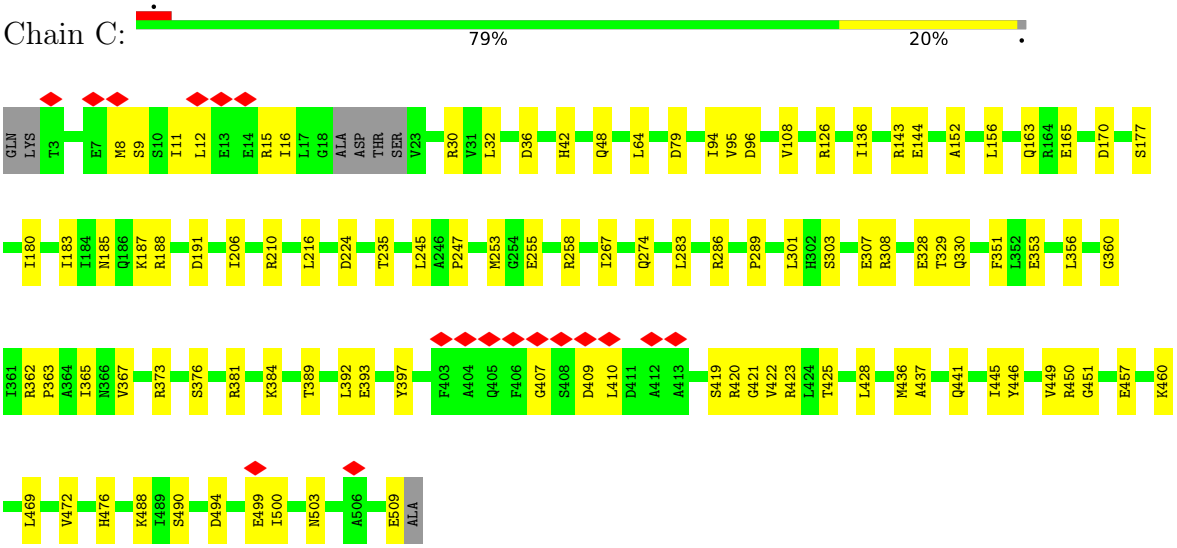


Mol	Chain	Residues	Atoms					AltConf
7	F	1	Total	C	N	O	P	0
			27	10	5	10	2	
7	D	1	Total	C	N	O	P	0
			27	10	5	10	2	

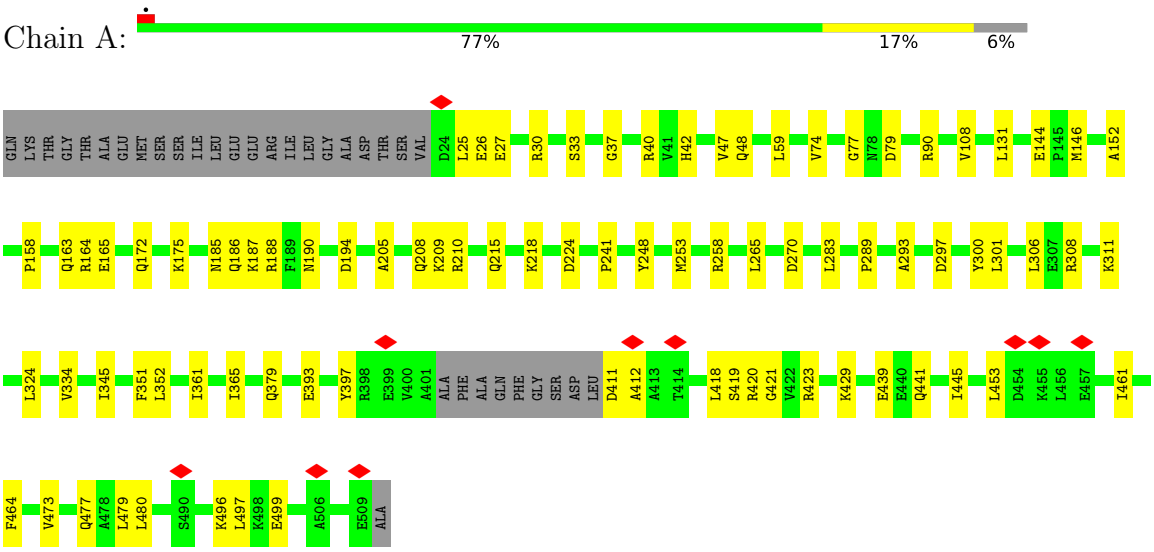
3 Residue-property plots (i)

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 subunit alpha, mitochondrial

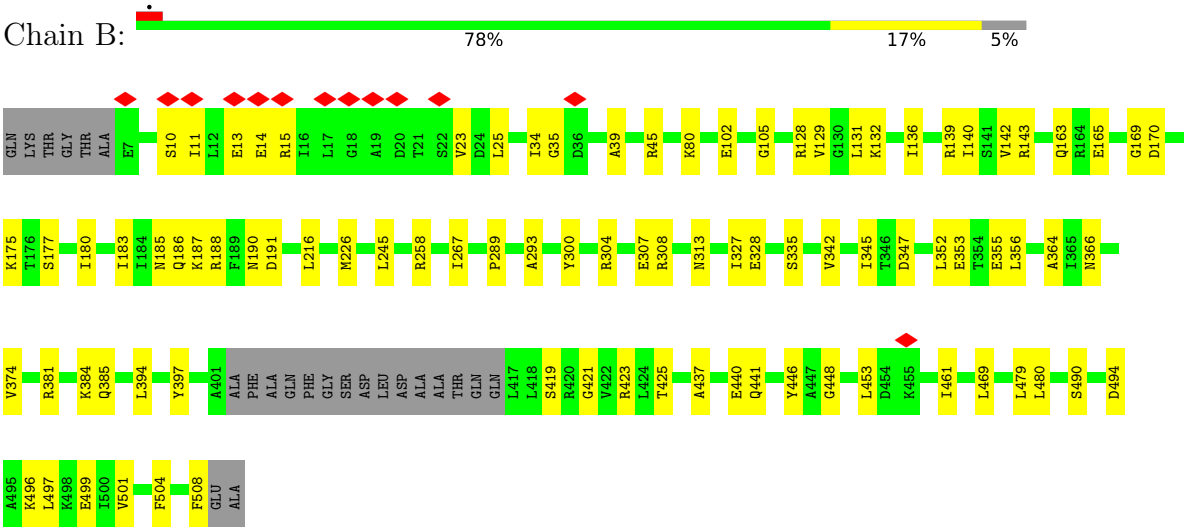


• Molecule 1: ATP synthase subunit alpha, mitochondrial



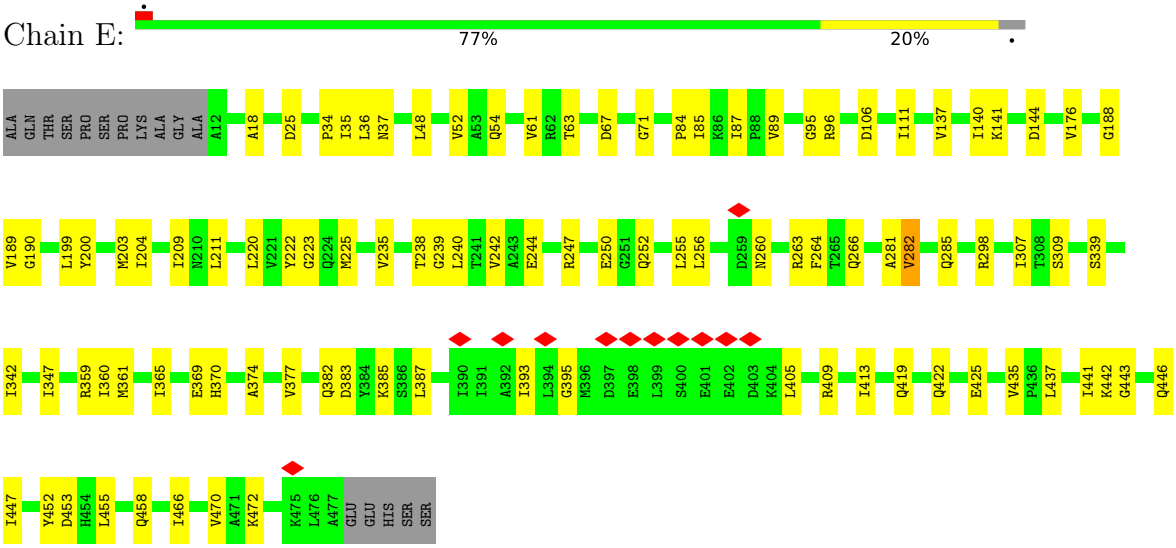
• Molecule 1: ATP synthase subunit alpha, mitochondrial

Chain B:



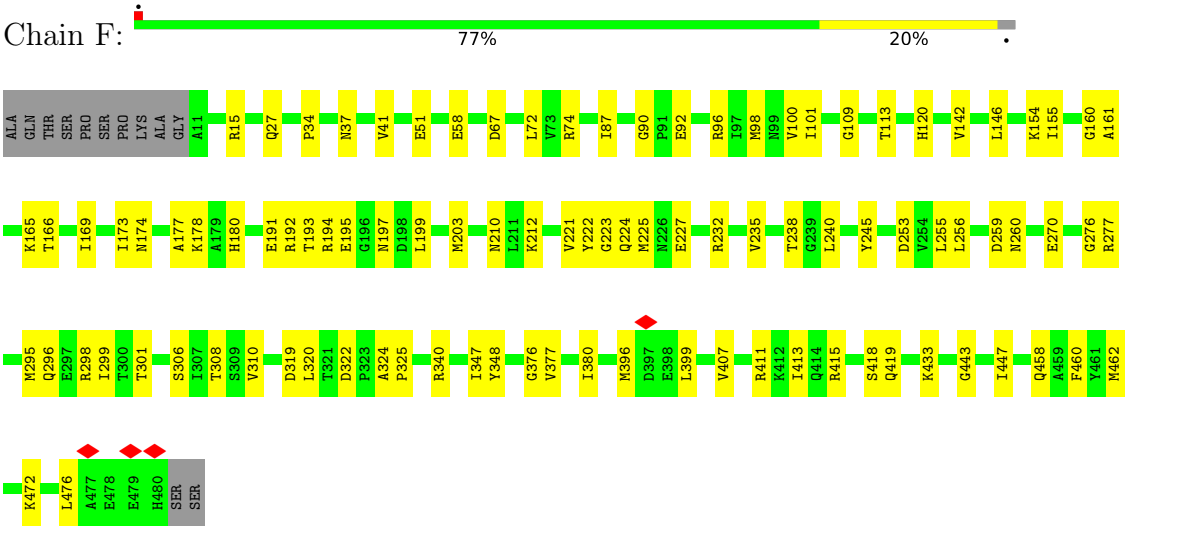
• Molecule 2: ATP synthase subunit beta, mitochondrial

Chain E:



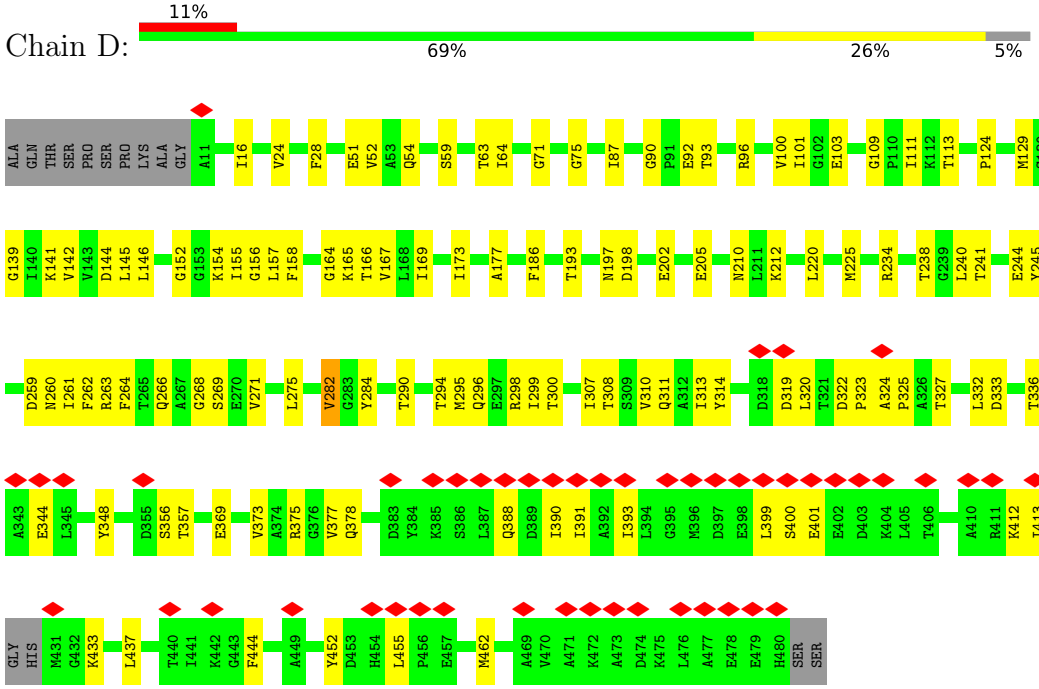
• Molecule 2: ATP synthase subunit beta, mitochondrial

Chain F:



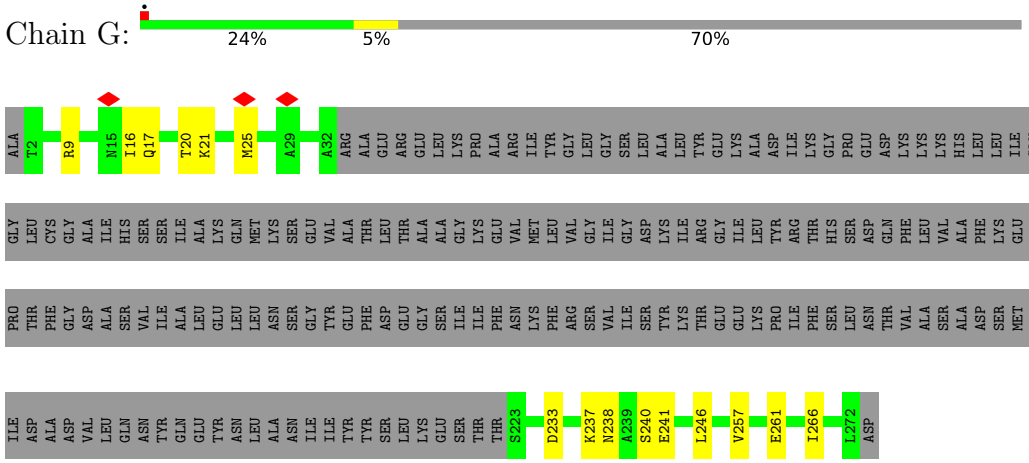
• Molecule 2: ATP synthase subunit beta, mitochondrial

Chain D:



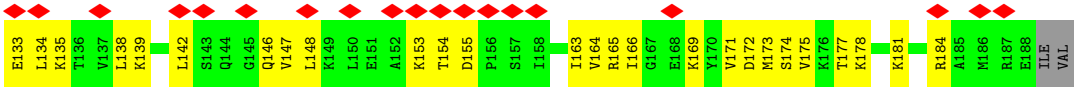
• Molecule 3: ATP synthase subunit gamma, mitochondrial

Chain G:



• Molecule 4: ATP synthase subunit O, mitochondrial

Chain O:


E133
L134
K135
T136
V137
L138
K139
L142
S143
Q144
G145
Q146
V147
L148
K149
L150
E151
A152
K153
T154
D155
P156
S157
I158
I163
V164
R165
I166
G167
E168
K169
Y170
V171
D172
M173
S174
V175
K176
T177
K178
K181
R184
A185
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	37480	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	0.785	Depositor
Minimum map value	-0.433	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.023	Depositor
Recommended contour level	0.1	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, ATP, ADP

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	A	0.26	0/3687	0.41	0/4973
1	B	0.28	0/3762	0.41	0/5073
1	C	0.28	0/3880	0.41	0/5233
2	D	0.27	0/3537	0.43	0/4795
2	E	0.31	0/3580	0.48	0/4857
2	F	0.29	0/3621	0.40	0/4911
3	G	0.21	0/623	0.35	0/829
4	O	0.20	0/1453	0.43	0/1958
All	All	0.28	0/24143	0.42	0/32629

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

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	A	3639	0	3741	54	0
1	B	3714	0	3825	63	0
1	C	3830	0	3928	70	0
2	D	3482	0	3536	89	0
2	E	3522	0	3572	66	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	F	3562	0	3609	60	0
3	G	623	0	677	10	0
4	O	1437	0	1536	51	0
5	A	31	0	12	0	0
5	B	31	0	12	2	0
5	C	31	0	12	1	0
6	A	1	0	0	0	0
6	B	1	0	0	0	0
6	C	1	0	0	0	0
6	D	1	0	0	0	0
6	F	1	0	0	0	0
7	D	27	0	12	6	0
7	F	27	0	12	0	0
All	All	23961	0	24484	429	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 9.

All (429) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:440:GLU:HB3	1:B:469:LEU:HD11	1.58	0.83
1:C:8:MET:HE3	2:F:58:GLU:H	1.48	0.77
4:O:135:LYS:HE2	4:O:139:LYS:HD2	1.66	0.76
2:F:174:ASN:O	2:F:178:LYS:HB2	1.87	0.76
2:D:388:GLN:HA	2:D:391:ILE:HG12	1.71	0.72
1:C:95:VAL:HG11	1:C:245:LEU:HD21	1.70	0.72
1:A:209:LYS:NZ	2:D:333:ASP:OD1	2.22	0.72
1:C:136:ILE:O	2:D:197:ASN:ND2	2.23	0.71
1:C:187:LYS:HE2	1:C:224:ASP:HB3	1.72	0.71
2:E:203:MET:HE1	2:E:220:LEU:HD21	1.72	0.70
3:G:16:ILE:HG12	2:D:393:ILE:HD11	1.73	0.70
2:D:141:LYS:O	2:D:145:LEU:HB3	1.91	0.70
2:E:96:ARG:HD3	2:E:111:ILE:HG12	1.73	0.69
1:A:30:ARG:HB3	4:O:59:TYR:CZ	2.27	0.69
2:D:152:GLY:H	2:D:308:THR:HB	1.57	0.69
2:D:158:PHE:HB2	2:D:337:VAL:HG22	1.75	0.69
1:C:185:ASN:OD1	1:C:188:ARG:NH1	2.26	0.69
2:F:227:GLU:O	2:F:232:ARG:NH1	2.26	0.69
2:E:244:GLU:OE2	2:E:298:ARG:NH1	2.26	0.68
2:F:161:ALA:O	2:F:340:ARG:NH1	2.27	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:261:ILE:HD11	2:D:295:MET:HE1	1.74	0.68
1:A:175:LYS:HG2	1:A:352:LEU:HD12	1.76	0.67
2:E:190:GLY:O	2:E:263:ARG:NE	2.27	0.67
1:A:205:ALA:HB1	1:A:208:GLN:HG3	1.75	0.67
1:C:286:ARG:NH2	2:F:276:GLY:O	2.28	0.66
2:F:210:ASN:ND2	2:F:212:LYS:O	2.28	0.66
4:O:139:LYS:HG2	4:O:148:LEU:HD13	1.77	0.66
2:D:164:GLY:HA2	7:D:501:ADP:H5'1	1.77	0.66
1:B:23:VAL:HG21	4:O:173:MET:HG2	1.78	0.66
1:C:499:GLU:OE2	1:C:503:ASN:ND2	2.27	0.65
1:A:393:GLU:OE2	1:A:420:ARG:NH1	2.30	0.65
4:O:115:GLU:HG3	4:O:147:VAL:HB	1.77	0.65
1:C:247:PRO:HG2	1:C:274:GLN:HG3	1.79	0.65
1:A:59:LEU:HD22	1:A:77:GLY:HA3	1.77	0.65
2:F:96:ARG:NH2	2:F:109:GLY:O	2.29	0.65
1:B:136:ILE:O	2:F:197:ASN:ND2	2.31	0.64
2:F:160:GLY:O	2:F:165:LYS:NZ	2.30	0.64
2:D:261:ILE:O	2:D:264:PHE:HB3	1.97	0.64
2:D:167:VAL:HG21	7:D:501:ADP:H8	1.63	0.64
2:E:250:GLU:OE1	2:E:252:GLN:NE2	2.31	0.64
2:D:210:ASN:ND2	2:D:212:LYS:O	2.30	0.63
1:A:163:GLN:NE2	1:A:165:GLU:OE2	2.31	0.63
2:E:141:LYS:H	2:E:419:GLN:HE22	1.44	0.63
1:C:457:GLU:OE2	1:C:460:LYS:NZ	2.28	0.63
4:O:142:LEU:HD23	4:O:146:GLN:HB3	1.81	0.63
1:B:185:ASN:OD1	1:B:188:ARG:NH1	2.32	0.63
1:C:36:ASP:OD1	2:F:277:ARG:NH2	2.32	0.62
4:O:110:SER:HA	4:O:113:ARG:HH21	1.65	0.62
4:O:125:PRO:HA	4:O:154:THR:HG21	1.81	0.62
2:D:375:ARG:NH1	2:D:378:GLN:OE1	2.32	0.62
1:C:490:SER:O	1:C:494:ASP:N	2.32	0.62
2:D:244:GLU:OE2	2:D:298:ARG:NH2	2.28	0.62
1:C:187:LYS:NZ	1:C:191:ASP:OD2	2.33	0.62
4:O:7:PRO:HG3	4:O:108:MET:HE3	1.80	0.62
2:D:413:ILE:HG23	2:D:444:PHE:HE1	1.65	0.62
1:A:37:GLY:N	1:A:79:ASP:OD2	2.34	0.61
1:A:423:ARG:HG3	1:A:461:ILE:HD11	1.82	0.61
2:D:156:GLY:HA3	2:D:332:LEU:HD13	1.82	0.61
4:O:35:VAL:HG11	4:O:81:LEU:HD23	1.83	0.61
2:D:369:GLU:O	2:D:373:VAL:HG23	2.00	0.60
2:F:92:GLU:HB2	2:F:113:THR:HG22	1.81	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:215:GLN:HG2	2:D:131:VAL:HG13	1.82	0.60
1:A:258:ARG:NH1	1:A:308:ARG:O	2.34	0.60
2:E:455:LEU:HD21	2:E:470:VAL:HG12	1.84	0.60
2:D:92:GLU:HB2	2:D:113:THR:HG22	1.83	0.59
1:C:9:SER:N	2:F:58:GLU:OE1	2.33	0.59
1:A:186:GLN:O	1:A:190:ASN:ND2	2.28	0.59
2:F:301:THR:HG23	2:F:306:SER:HA	1.83	0.59
2:F:411:ARG:O	2:F:415:ARG:HG2	2.02	0.59
2:D:166:THR:O	2:D:169:ILE:HG22	2.03	0.59
1:A:187:LYS:HE2	1:A:224:ASP:HB3	1.85	0.59
1:B:258:ARG:NH1	1:B:308:ARG:O	2.35	0.59
1:B:10:SER:O	1:B:15:ARG:NH2	2.35	0.58
1:B:163:GLN:NE2	1:B:165:GLU:OE2	2.36	0.58
2:D:247:ARG:HD3	2:D:307:ILE:HG13	1.84	0.58
1:A:185:ASN:OD1	1:A:188:ARG:NH1	2.37	0.58
2:E:141:LYS:H	2:E:419:GLN:NE2	2.00	0.58
2:D:262:PHE:HB3	2:D:314:TYR:HB3	1.86	0.58
1:C:156:LEU:HD13	1:C:367:VAL:HG11	1.85	0.58
2:F:458:GLN:O	2:F:472:LYS:NZ	2.32	0.58
2:D:157:LEU:N	2:D:311:GLN:O	2.37	0.57
1:A:218:LYS:HD3	2:D:131:VAL:HG11	1.86	0.57
3:G:17:GLN:NE2	3:G:240:SER:OG	2.37	0.57
2:F:15:ARG:NH2	2:F:27:GLN:OE1	2.31	0.57
2:F:347:ILE:HG23	2:F:418:SER:HB3	1.85	0.57
4:O:172:ASP:OD2	4:O:177:THR:HG21	2.04	0.57
2:D:268:GLY:O	2:D:271:VAL:HG22	2.05	0.57
2:E:247:ARG:HD3	2:E:307:ILE:HG13	1.87	0.57
1:A:479:LEU:HD21	1:A:497:LEU:HG	1.86	0.57
2:F:51:GLU:OE2	2:F:120:HIS:NE2	2.36	0.57
4:O:39:LEU:HD13	4:O:102:VAL:HG13	1.86	0.57
2:D:154:LYS:NZ	2:D:296:GLN:O	2.28	0.57
1:B:140:ILE:HG12	1:B:143:ARG:HH12	1.70	0.57
1:C:96:ASP:OD1	1:C:126:ARG:NE	2.31	0.56
1:B:175:LYS:HG2	1:B:352:LEU:HD12	1.87	0.56
3:G:21:LYS:O	3:G:25:MET:HG2	2.04	0.56
4:O:8:PRO:HG2	4:O:111:VAL:HG23	1.86	0.56
1:B:186:GLN:O	1:B:190:ASN:ND2	2.28	0.56
2:E:84:PRO:O	2:E:85:ILE:HG13	2.06	0.56
2:D:139:GLY:N	2:D:144:ASP:OD2	2.34	0.56
2:D:255:LEU:HD23	2:D:308:THR:HG23	1.85	0.56
1:A:439:GLU:HG3	1:A:480:LEU:HB3	1.88	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:240:LEU:HD21	2:E:298:ARG:HB2	1.87	0.56
1:C:163:GLN:NE2	1:C:165:GLU:OE2	2.39	0.56
1:A:473:VAL:O	1:A:477:GLN:NE2	2.37	0.56
2:D:165:LYS:H	7:D:501:ADP:PB	2.27	0.56
1:C:177:SER:OG	5:C:601:ATP:O1A	2.24	0.56
1:B:496:LYS:O	1:B:499:GLU:HG3	2.05	0.56
1:B:25:LEU:HG	4:O:169:LYS:HE3	1.88	0.55
1:C:397:TYR:CG	1:C:421:GLY:HA3	2.41	0.55
2:E:393:ILE:HG23	2:E:395:GLY:H	1.72	0.55
2:E:263:ARG:HA	2:E:266:GLN:HG2	1.88	0.55
4:O:18:TYR:OH	4:O:92:ASN:ND2	2.38	0.55
2:F:34:PRO:HD2	2:F:37:ASN:ND2	2.22	0.55
2:D:420:PRO:HG2	2:D:433:LYS:HG2	1.89	0.55
1:B:307:GLU:OE2	2:F:193:THR:OG1	2.22	0.54
1:B:13:GLU:HG3	1:B:14:GLU:H	1.72	0.54
1:C:152:ALA:HB3	1:C:365:ILE:HD12	1.88	0.54
2:F:232:ARG:NH2	2:F:270:GLU:OE1	2.41	0.54
2:D:100:VAL:HG13	2:D:101:ILE:HG23	1.88	0.54
1:C:328:GLU:OE1	1:C:330:GLN:NE2	2.40	0.54
2:F:191:GLU:HG3	2:F:192:ARG:H	1.73	0.54
1:B:347:ASP:OD1	2:F:194:ARG:NE	2.40	0.54
1:C:362:ARG:HA	1:C:363:PRO:C	2.33	0.54
1:B:187:LYS:NZ	1:B:191:ASP:OD2	2.41	0.54
1:A:283:LEU:HD11	1:A:293:ALA:HB1	1.89	0.53
2:E:422:GLN:O	2:E:425:GLU:HB2	2.09	0.53
2:E:442:LYS:HE3	2:E:466:ILE:HD11	1.89	0.53
2:D:101:ILE:HG13	2:D:103:GLU:HG3	1.90	0.53
1:C:183:ILE:HD11	1:C:267:ILE:HD13	1.90	0.53
1:C:381:ARG:HB3	1:C:488:LYS:HB3	1.91	0.53
1:B:300:TYR:O	1:B:304:ARG:HG2	2.08	0.53
2:F:348:TYR:HB2	2:F:462:MET:HE1	1.90	0.53
1:C:143:ARG:HG3	1:C:144:GLU:HG3	1.91	0.53
2:E:87:ILE:HG21	2:E:238:THR:HG23	1.91	0.53
2:D:377:VAL:HG13	2:D:413:ILE:HG21	1.91	0.53
4:O:129:ALA:O	4:O:133:GLU:HG3	2.10	0.52
2:D:369:GLU:CD	2:D:369:GLU:H	2.16	0.52
2:E:458:GLN:O	2:E:472:LYS:NZ	2.43	0.52
2:E:18:ALA:HB3	2:E:25:ASP:HB2	1.92	0.52
2:E:443:GLY:HA2	2:E:466:ILE:HD13	1.92	0.52
1:A:361:ILE:HD13	1:A:429:LYS:HE2	1.91	0.52
2:E:383:ASP:O	2:E:387:LEU:HG	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:141:LYS:O	2:D:145:LEU:CB	2.56	0.52
1:A:108:VAL:HG21	1:A:253:MET:HE1	1.91	0.51
1:A:283:LEU:HD21	1:A:289:PRO:HB3	1.91	0.51
4:O:65:LYS:HD3	4:O:95:LEU:HD13	1.91	0.51
2:D:186:PHE:HB3	2:D:220:LEU:HD23	1.90	0.51
1:C:373:ARG:NH2	7:D:501:ADP:O3B	2.43	0.51
1:B:139:ARG:NH2	1:B:307:GLU:O	2.41	0.51
1:C:420:ARG:HH22	1:C:451:GLY:HA2	1.75	0.51
1:A:419:SER:O	1:A:423:ARG:HD3	2.11	0.51
2:D:400:SER:OG	2:D:401:GLU:N	2.43	0.51
4:O:69:LEU:HD12	4:O:72:ILE:HD11	1.91	0.51
4:O:181:LYS:HG2	4:O:184:ARG:HH21	1.75	0.51
2:F:322:ASP:O	2:F:325:PRO:HD2	2.11	0.51
1:B:353:GLU:HG3	1:B:366:ASN:HB2	1.93	0.51
1:B:170:ASP:O	1:B:175:LYS:NZ	2.43	0.51
2:D:169:ILE:HD13	2:D:259:ASP:HB2	1.92	0.51
1:C:210:ARG:HG3	1:C:235:THR:HG21	1.93	0.51
2:D:340:ARG:O	2:D:344:GLU:HG3	2.10	0.50
1:B:180:ILE:HD11	1:B:216:LEU:HD11	1.93	0.50
2:D:54:GLN:NE2	2:D:275:LEU:O	2.37	0.50
1:B:397:TYR:CG	1:B:421:GLY:HA3	2.47	0.50
2:E:339:SER:HB2	2:E:342:ILE:HG12	1.93	0.50
2:D:111:ILE:HG22	2:D:113:THR:HG23	1.94	0.50
1:C:255:GLU:OE1	1:C:258:ARG:NH2	2.45	0.50
1:B:453:LEU:HD12	1:B:461:ILE:HG12	1.93	0.50
2:F:240:LEU:HD21	2:F:298:ARG:HB2	1.94	0.50
1:B:448:GLY:HA2	1:B:453:LEU:HD22	1.93	0.50
2:F:87:ILE:HD13	2:F:238:THR:HG23	1.93	0.50
1:B:13:GLU:HG3	1:B:14:GLU:N	2.27	0.50
3:G:9:ARG:HE	3:G:246:LEU:HD11	1.77	0.50
1:B:446:TYR:OH	1:B:494:ASP:OD1	2.30	0.49
2:E:281:ALA:O	2:E:282:VAL:HB	2.11	0.49
2:E:405:LEU:O	2:E:409:ARG:HG2	2.12	0.49
1:A:210:ARG:NH1	2:D:124:PRO:O	2.44	0.49
1:B:490:SER:O	1:B:494:ASP:N	2.42	0.49
1:C:420:ARG:NH2	1:C:451:GLY:HA2	2.27	0.49
1:C:15:ARG:HG3	4:O:88:LEU:HD11	1.95	0.49
1:C:170:ASP:OD1	1:C:329:THR:OG1	2.30	0.49
2:F:173:ILE:O	2:F:177:ALA:HB3	2.13	0.49
2:D:324:ALA:HB3	2:D:325:PRO:HD3	1.95	0.49
2:F:443:GLY:O	2:F:447:ILE:HG13	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:335:SER:O	1:B:335:SER:OG	2.27	0.49
1:B:381:ARG:HG3	1:B:385:GLN:HE21	1.77	0.48
2:F:199:LEU:HD11	2:F:203:MET:HE3	1.95	0.48
4:O:69:LEU:HG	4:O:86:ILE:HD12	1.95	0.48
4:O:76:GLU:C	4:O:77:ARG:HD3	2.38	0.48
2:F:41:VAL:HG11	2:F:72:LEU:HD21	1.95	0.48
1:B:353:GLU:OE2	1:B:366:ASN:ND2	2.41	0.48
2:F:419:GLN:NE2	2:F:433:LYS:O	2.42	0.48
3:G:233:ASP:O	3:G:237:LYS:HG2	2.14	0.48
2:D:348:TYR:HB2	2:D:462:MET:HE1	1.95	0.48
1:B:289:PRO:HB2	1:B:293:ALA:HA	1.95	0.48
4:O:39:LEU:HD21	4:O:105:PHE:CD1	2.48	0.48
1:A:164:ARG:NH1	1:A:306:LEU:O	2.40	0.48
2:E:360:ILE:HB	2:E:365:ILE:HG21	1.95	0.48
4:O:121:THR:HA	4:O:153:LYS:O	2.14	0.48
1:C:8:MET:HE3	2:F:58:GLU:N	2.23	0.48
2:E:52:VAL:HA	2:E:63:THR:HG22	1.96	0.48
1:A:30:ARG:HB3	4:O:59:TYR:OH	2.14	0.48
1:C:258:ARG:NH1	1:C:308:ARG:O	2.47	0.47
1:A:172:GLN:HG2	2:D:357:THR:HG21	1.95	0.47
1:B:105:GLY:HA2	1:B:226:MET:O	2.15	0.47
1:A:74:VAL:HB	1:A:241:PRO:HG3	1.97	0.47
3:G:237:LYS:NZ	3:G:238:ASN:OD1	2.47	0.47
1:B:356:LEU:HB2	1:B:364:ALA:HB1	1.95	0.47
2:E:141:LYS:HG3	2:E:435:VAL:HG11	1.95	0.47
2:E:370:HIS:CD2	2:E:441:ILE:HD11	2.50	0.47
2:F:377:VAL:HG13	2:F:413:ILE:HG21	1.96	0.47
1:C:389:THR:O	1:C:393:GLU:HG3	2.15	0.47
2:E:377:VAL:HG13	2:E:413:ILE:HG21	1.97	0.47
4:O:36:GLU:HB2	4:O:109:MET:SD	2.54	0.47
4:O:69:LEU:HD21	4:O:86:ILE:HB	1.96	0.47
1:C:437:ALA:O	1:C:441:GLN:HG3	2.15	0.47
1:A:25:LEU:HD21	1:A:42:HIS:CE1	2.50	0.47
1:B:381:ARG:NH1	1:B:384:LYS:HD2	2.30	0.47
1:B:421:GLY:O	1:B:425:THR:HG23	2.15	0.47
2:F:155:ILE:HD12	2:F:310:VAL:HG22	1.97	0.47
4:O:70:ASN:HA	4:O:73:THR:HG22	1.96	0.47
1:C:303:SER:OG	2:D:225:MET:HB3	2.15	0.47
4:O:171:VAL:HG12	4:O:173:MET:HG3	1.97	0.47
2:D:52:VAL:HA	2:D:63:THR:HG22	1.97	0.47
2:E:204:ILE:HD13	2:E:211:LEU:HD11	1.95	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:322:ASP:HB3	2:D:325:PRO:HD2	1.96	0.46
4:O:104:ALA:O	4:O:108:MET:HG3	2.15	0.46
1:B:132:LYS:NZ	2:F:67:ASP:OD2	2.49	0.46
2:F:98:MET:HG3	2:F:221:VAL:HG22	1.98	0.46
4:O:69:LEU:HD22	4:O:95:LEU:HD11	1.98	0.46
4:O:118:CYS:HA	4:O:165:ARG:O	2.16	0.46
2:E:256:LEU:O	2:E:309:SER:HA	2.16	0.46
1:C:509:GLU:OE1	1:C:509:GLU:N	2.48	0.46
2:D:173:ILE:O	2:D:177:ALA:HB3	2.15	0.46
2:E:200:TYR:CZ	2:E:204:ILE:HD11	2.50	0.46
2:F:100:VAL:HG13	2:F:101:ILE:HG23	1.98	0.46
4:O:127:GLU:H	4:O:127:GLU:CD	2.23	0.46
2:D:87:ILE:HD13	2:D:238:THR:HG23	1.97	0.46
1:C:79:ASP:OD1	1:C:79:ASP:N	2.49	0.46
2:D:137:VAL:HG11	2:D:437:LEU:HB2	1.98	0.46
1:C:48:GLN:HB3	2:D:71:GLY:HA2	1.98	0.46
3:G:237:LYS:O	3:G:241:GLU:HG3	2.16	0.46
2:D:390:ILE:HG22	2:D:399:LEU:HD11	1.98	0.45
1:C:392:LEU:HD12	1:C:393:GLU:N	2.31	0.45
4:O:123:ALA:HA	4:O:155:ASP:O	2.16	0.45
2:D:412:LYS:NZ	2:D:452:TYR:O	2.50	0.45
2:E:140:ILE:HA	2:E:419:GLN:HE22	1.80	0.45
4:O:15:GLU:OE2	4:O:97:ASN:HB3	2.16	0.45
4:O:175:VAL:HA	4:O:178:LYS:HE3	1.98	0.45
1:C:436:MET:HE2	1:C:441:GLN:HG2	1.99	0.45
1:C:469:LEU:HA	1:C:472:VAL:HG22	1.98	0.45
4:O:127:GLU:OE1	4:O:127:GLU:N	2.39	0.45
2:D:87:ILE:HD11	2:D:241:THR:HB	1.98	0.45
2:D:164:GLY:CA	7:D:501:ADP:H5'1	2.44	0.45
1:C:389:THR:HG23	1:C:449:VAL:HG21	1.97	0.45
1:A:48:GLN:HB3	2:E:71:GLY:HA2	1.99	0.45
2:E:34:PRO:HD2	2:E:37:ASN:HD22	1.82	0.45
2:F:324:ALA:HB3	2:F:325:PRO:HD3	1.99	0.45
2:D:240:LEU:HD21	2:D:299:ILE:HG12	1.98	0.45
2:E:382:GLN:O	2:E:385:LYS:HG2	2.17	0.45
1:C:389:THR:HA	1:C:392:LEU:HD21	1.98	0.44
1:B:497:LEU:O	1:B:501:VAL:HG12	2.17	0.44
2:E:443:GLY:O	2:E:447:ILE:HG13	2.17	0.44
4:O:163:ILE:HG12	4:O:172:ASP:OD1	2.18	0.44
1:C:419:SER:O	1:C:423:ARG:HD3	2.17	0.44
1:B:169:GLY:O	1:B:328:GLU:HA	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:52:VAL:HG13	2:E:61:VAL:HG13	1.99	0.44
2:F:180:HIS:HE1	2:F:253:ASP:O	2.01	0.44
2:D:261:ILE:HG22	2:D:313:ILE:HG12	1.99	0.44
1:A:345:ILE:HA	2:E:225:MET:HE1	2.00	0.44
1:B:504:PHE:O	1:B:508:PHE:N	2.48	0.44
2:D:87:ILE:HG21	2:D:238:THR:HG23	1.98	0.44
2:D:240:LEU:HD21	2:D:299:ILE:CG1	2.47	0.44
1:B:142:VAL:HB	1:B:374:VAL:HG11	1.99	0.44
2:E:95:GLY:HA2	2:E:209:ILE:HG12	1.99	0.44
2:E:188:GLY:HA3	2:E:222:TYR:HD1	1.82	0.44
2:F:259:ASP:HA	2:F:260:ASN:HA	1.75	0.44
2:D:198:ASP:O	2:D:202:GLU:HG3	2.18	0.44
1:A:146:MET:HE3	1:A:146:MET:HB3	1.92	0.44
1:B:80:LYS:HA	2:E:35:ILE:HB	2.00	0.44
2:E:409:ARG:NH2	2:E:453:ASP:HB3	2.32	0.44
2:D:263:ARG:HA	2:D:266:GLN:HB3	1.99	0.44
1:A:144:GLU:OE2	1:A:311:LYS:HE2	2.18	0.44
1:B:34:ILE:HD13	1:B:39:ALA:HB2	1.99	0.44
1:B:353:GLU:CD	1:B:366:ASN:HD22	2.24	0.44
2:E:199:LEU:HD23	2:E:222:TYR:OH	2.17	0.44
1:C:206:ILE:HD11	1:C:247:PRO:HG3	2.00	0.44
1:C:446:TYR:CE1	1:C:450:ARG:HG3	2.53	0.44
2:E:37:ASN:HA	2:E:84:PRO:HA	1.99	0.44
2:E:361:MET:HE1	2:E:374:ALA:HB3	1.98	0.44
2:F:396:MET:HG3	2:F:399:LEU:HD12	2.00	0.44
2:F:154:LYS:NZ	2:F:296:GLN:O	2.51	0.44
4:O:89:LEU:HB3	4:O:95:LEU:HA	2.00	0.44
1:A:248:TYR:OH	1:A:301:LEU:O	2.31	0.44
1:B:381:ARG:O	1:B:385:GLN:HG3	2.17	0.44
2:D:259:ASP:HA	2:D:260:ASN:HA	1.69	0.44
1:C:360:GLY:HA2	1:C:362:ARG:NH2	2.33	0.43
1:A:265:LEU:HD11	1:A:324:LEU:HG	1.99	0.43
1:A:397:TYR:CG	1:A:421:GLY:HA3	2.53	0.43
1:A:411:ASP:OD1	1:A:412:ALA:N	2.51	0.43
2:D:290:THR:O	2:D:294:THR:HG23	2.18	0.43
2:E:442:LYS:HB3	2:E:442:LYS:HE2	1.64	0.43
1:C:64:LEU:HD23	1:C:64:LEU:HA	1.84	0.43
1:A:453:LEU:HD23	1:A:461:ILE:HG13	2.00	0.43
2:D:64:ILE:HD11	2:D:275:LEU:HD11	2.00	0.43
2:D:96:ARG:NH2	2:D:109:GLY:O	2.30	0.43
2:F:396:MET:HE2	2:F:407:VAL:HG11	1.99	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:334:VAL:HG11	1:A:351:PHE:CE1	2.54	0.43
1:B:11:ILE:HA	1:B:15:ARG:HH21	1.84	0.43
2:E:89:VAL:HG23	2:E:242:VAL:HG23	2.00	0.43
2:F:166:THR:HA	2:F:169:ILE:HG22	2.00	0.43
2:D:51:GLU:CD	2:D:234:ARG:HE	2.27	0.43
2:D:433:LYS:HA	2:D:433:LYS:HD3	1.81	0.43
1:C:180:ILE:HD11	1:C:216:LEU:HD11	2.00	0.43
1:C:476:HIS:CE1	1:C:500:ILE:HG12	2.54	0.43
1:A:152:ALA:HB3	1:A:365:ILE:HD12	2.00	0.43
1:A:441:GLN:O	1:A:445:ILE:HG12	2.18	0.43
1:B:183:ILE:HD11	1:B:267:ILE:HD13	2.01	0.43
1:B:479:LEU:HD23	1:B:480:LEU:HD23	2.00	0.43
2:F:222:TYR:HD2	2:F:224:GLN:NE2	2.17	0.43
2:F:472:LYS:O	2:F:476:LEU:HG	2.19	0.43
4:O:164:VAL:HG12	4:O:171:VAL:HB	2.01	0.43
2:D:452:TYR:HB3	2:D:455:LEU:HD12	2.00	0.43
1:C:108:VAL:HG21	1:C:253:MET:HE1	2.01	0.42
1:C:445:ILE:O	1:C:449:VAL:HG12	2.19	0.42
1:A:26:GLU:HG3	1:A:27:GLU:OE1	2.18	0.42
1:B:11:ILE:O	1:B:15:ARG:HB2	2.19	0.42
1:B:140:ILE:HB	1:B:313:ASN:HB3	2.01	0.42
1:B:419:SER:O	1:B:423:ARG:HD3	2.19	0.42
2:E:106:ASP:OD1	2:E:106:ASP:N	2.51	0.42
1:C:376:SER:HB2	1:C:384:LYS:HG3	2.01	0.42
2:D:412:LYS:NZ	2:D:455:LEU:O	2.48	0.42
1:B:177:SER:OG	5:B:601:ATP:O1A	2.33	0.42
2:E:176:VAL:HG21	2:E:255:LEU:HD22	2.00	0.42
2:E:446:GLN:NE2	2:E:452:TYR:OH	2.53	0.42
2:F:376:GLY:O	2:F:380:ILE:HG12	2.19	0.42
3:G:257:VAL:O	3:G:261:GLU:HG2	2.19	0.42
2:E:342:ILE:HG22	2:E:347:ILE:HB	2.00	0.42
2:D:167:VAL:HG21	7:D:501:ADP:C8	2.52	0.42
4:O:134:LEU:O	4:O:138:LEU:HG	2.20	0.42
1:A:131:LEU:HD23	1:A:131:LEU:HA	1.90	0.42
2:D:202:GLU:HA	2:D:205:GLU:HG3	2.02	0.42
1:A:297:ASP:HB2	1:A:300:TYR:HB3	2.01	0.42
2:F:192:ARG:HB2	2:F:195:GLU:HG3	2.02	0.42
2:F:319:ASP:O	2:F:320:LEU:HB2	2.18	0.42
1:C:407:GLY:HA2	1:C:410:LEU:HD13	2.02	0.42
2:F:256:LEU:HD23	2:F:299:ILE:HG23	2.01	0.42
3:G:266:ILE:HD13	3:G:266:ILE:HA	1.86	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:274:GLN:HE22	1:C:301:LEU:HD11	1.84	0.42
2:E:223:GLY:N	2:E:235:VAL:HG21	2.34	0.42
1:C:409:ASP:C	1:C:410:LEU:HD12	2.45	0.42
1:C:428:LEU:HD23	1:C:428:LEU:HA	1.85	0.42
1:A:194:ASP:OD2	1:A:194:ASP:N	2.52	0.42
2:F:142:VAL:HG23	2:F:146:LEU:HD12	2.02	0.42
2:D:323:PRO:O	2:D:327:THR:OG1	2.29	0.42
1:A:334:VAL:HG11	1:A:351:PHE:HE1	1.85	0.41
5:B:601:ATP:O2G	2:E:359:ARG:NH1	2.45	0.41
2:D:154:LYS:HE3	2:D:299:ILE:HB	2.01	0.41
4:O:7:PRO:HG3	4:O:108:MET:CE	2.50	0.41
4:O:89:LEU:HD13	4:O:98:THR:HB	2.02	0.41
2:D:16:ILE:O	2:D:75:GLY:N	2.48	0.41
2:D:142:VAL:HG23	2:D:146:LEU:HD12	2.02	0.41
1:C:12:LEU:HD22	4:O:88:LEU:HD23	2.01	0.41
1:A:47:VAL:HA	1:A:90:ARG:HD2	2.02	0.41
1:A:397:TYR:HE2	1:A:418:LEU:HD12	1.85	0.41
1:B:423:ARG:HG3	1:B:461:ILE:HD11	2.02	0.41
2:E:34:PRO:HD2	2:E:37:ASN:ND2	2.36	0.41
2:E:370:HIS:HD2	2:E:441:ILE:HD11	1.85	0.41
1:C:94:ILE:HG22	1:C:95:VAL:H	1.85	0.41
1:C:351:PHE:HE1	1:C:353:GLU:OE2	2.02	0.41
1:C:32:LEU:HD11	1:C:42:HIS:HB2	2.02	0.41
1:C:422:VAL:HA	1:C:425:THR:HG22	2.02	0.41
1:A:453:LEU:HD11	1:A:464:PHE:CE1	2.55	0.41
1:B:345:ILE:HA	2:F:225:MET:HE1	2.02	0.41
2:E:48:LEU:HD13	2:E:67:ASP:HB3	2.01	0.41
2:E:370:HIS:CD2	2:E:437:LEU:HD11	2.56	0.41
1:C:353:GLU:HB2	1:C:356:LEU:HB2	2.02	0.41
1:A:496:LYS:O	1:A:499:GLU:HG3	2.21	0.41
1:B:437:ALA:O	1:B:441:GLN:HG3	2.21	0.41
2:D:261:ILE:CG2	2:D:313:ILE:HG12	2.51	0.41
2:D:282:VAL:O	2:D:282:VAL:HG12	2.21	0.41
1:C:36:ASP:CG	2:F:277:ARG:HH21	2.28	0.41
1:C:11:ILE:HD11	1:C:30:ARG:HH12	1.85	0.41
1:A:158:PRO:HB3	1:A:379:GLN:HG3	2.03	0.41
1:B:394:LEU:HD23	1:B:394:LEU:HA	1.92	0.41
2:E:239:GLY:HA2	2:E:242:VAL:HG12	2.03	0.41
2:E:244:GLU:HA	2:E:307:ILE:HD11	2.03	0.41
2:E:260:ASN:HD22	2:E:263:ARG:NH1	2.19	0.41
2:E:285:GLN:H	2:E:285:GLN:HG2	1.52	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:90:GLY:HA2	2:F:245:TYR:CE1	2.55	0.41
2:F:255:LEU:HD23	2:F:308:THR:HB	2.01	0.41
2:D:90:GLY:HA2	2:D:245:TYR:CE1	2.56	0.41
2:D:155:ILE:HD12	2:D:310:VAL:HG22	2.03	0.41
1:C:16:ILE:HG23	4:O:88:LEU:HD22	2.03	0.41
1:C:283:LEU:HD21	1:C:289:PRO:HB3	2.03	0.41
1:A:42:HIS:HB3	4:O:59:TYR:OH	2.21	0.41
2:E:35:ILE:HG22	2:E:36:LEU:HG	2.02	0.41
2:E:189:VAL:HG11	2:E:264:PHE:HB2	2.02	0.41
2:E:369:GLU:HG2	2:E:370:HIS:H	1.85	0.41
2:D:269:SER:OG	2:D:284:TYR:HD1	2.04	0.41
1:B:129:VAL:HG21	1:B:245:LEU:HD11	2.03	0.40
2:E:137:VAL:HG13	2:E:144:ASP:OD2	2.22	0.40
1:B:35:GLY:HA3	2:E:54:GLN:HG3	2.03	0.40
1:B:45:ARG:NH1	2:F:74:ARG:HH21	2.19	0.40
2:F:223:GLY:N	2:F:235:VAL:HG21	2.36	0.40
2:F:295:MET:HB3	2:F:295:MET:HE3	1.84	0.40
2:F:415:ARG:HB2	2:F:460:PHE:HB2	2.02	0.40
4:O:138:LEU:HD22	4:O:166:ILE:HD11	2.02	0.40
2:D:24:VAL:O	2:D:63:THR:OG1	2.30	0.40
2:D:28:PHE:O	2:D:59:SER:HB3	2.22	0.40
2:D:93:THR:HG22	2:D:111:ILE:HG21	2.02	0.40
2:D:390:ILE:CG2	2:D:399:LEU:HD11	2.52	0.40
4:O:38:GLU:O	4:O:42:VAL:HG23	2.21	0.40
4:O:172:ASP:OD1	4:O:174:SER:OG	2.33	0.40
2:D:319:ASP:O	2:D:320:LEU:HB2	2.21	0.40
1:B:128:ARG:HB2	1:B:131:LEU:HG	2.02	0.40
1:B:327:ILE:HD11	1:B:342:VAL:HG21	2.03	0.40
1:B:355:GLU:OE2	1:B:355:GLU:N	2.51	0.40
1:B:448:GLY:HA2	1:B:453:LEU:CD2	2.51	0.40
4:O:69:LEU:HA	4:O:72:ILE:HG12	2.03	0.40
2:D:336:THR:HA	2:D:356:SER:HB2	2.04	0.40
1:C:307:GLU:OE2	2:D:193:THR:OG1	2.27	0.40
1:C:389:THR:O	1:C:392:LEU:HG	2.21	0.40
1:A:33:SER:HB3	1:A:40:ARG:HB2	2.04	0.40
1:A:270:ASP:OD1	1:A:270:ASP:N	2.54	0.40
1:B:102:GLU:OE1	1:B:102:GLU:N	2.43	0.40
3:G:16:ILE:O	3:G:20:THR:HG23	2.22	0.40
4:O:108:MET:O	4:O:111:VAL:HG22	2.22	0.40
2:D:129:MET:HE1	2:D:300:THR:HG21	2.03	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	A	473/510 (93%)	454 (96%)	19 (4%)	0	100	100
1	B	483/510 (95%)	468 (97%)	15 (3%)	0	100	100
1	C	499/510 (98%)	483 (97%)	16 (3%)	0	100	100
2	D	456/482 (95%)	428 (94%)	27 (6%)	1 (0%)	44	71
2	E	464/482 (96%)	432 (93%)	31 (7%)	1 (0%)	44	71
2	F	468/482 (97%)	446 (95%)	22 (5%)	0	100	100
3	G	77/273 (28%)	77 (100%)	0	0	100	100
4	O	185/190 (97%)	182 (98%)	3 (2%)	0	100	100
All	All	3105/3439 (90%)	2970 (96%)	133 (4%)	2 (0%)	50	76

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	E	282	VAL
2	D	282	VAL

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	A	387/412 (94%)	387 (100%)	0	100	100
1	B	397/412 (96%)	397 (100%)	0	100	100
1	C	407/412 (99%)	407 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	D	371/388 (96%)	371 (100%)	0	100	100
2	E	374/388 (96%)	374 (100%)	0	100	100
2	F	379/388 (98%)	379 (100%)	0	100	100
3	G	68/229 (30%)	68 (100%)	0	100	100
4	O	162/165 (98%)	162 (100%)	0	100	100
All	All	2545/2794 (91%)	2545 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (38) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	302	HIS
1	C	330	GLN
1	C	396	GLN
1	C	405	GLN
1	C	416	GLN
1	C	466	ASN
1	C	477	GLN
1	B	48	GLN
1	B	78	ASN
1	B	172	GLN
1	B	208	GLN
1	B	302	HIS
1	B	385	GLN
1	B	432	GLN
1	B	477	GLN
2	E	133	GLN
2	E	210	ASN
2	E	260	ASN
2	E	296	GLN
2	E	311	GLN
2	E	370	HIS
2	E	382	GLN
2	E	419	GLN
2	E	446	GLN
2	E	458	GLN
2	F	180	HIS
2	F	266	GLN
3	G	17	GLN

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Mol	Chain	Res	Type
4	O	70	ASN
4	O	84	ASN
4	O	92	ASN
4	O	112	HIS
4	O	146	GLN
2	D	133	GLN
2	D	201	HIS
2	D	296	GLN
2	D	311	GLN
2	D	370	HIS

5.3.3 RNA

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates

There are no oligosaccharides in this entry.

5.6 Ligand geometry

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	$\#\|Z\| > 2$
7	ADP	F	501	6	24,29,29	2.36	8 (33%)	29,45,45	1.64	6 (20%)
7	ADP	D	501	-	24,29,29	0.68	0	29,45,45	0.93	2 (6%)
5	ATP	B	601	6	26,33,33	2.20	6 (23%)	31,52,52	1.79	7 (22%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
5	ATP	A	601	6	26,33,33	2.25	6 (23%)	31,52,52	1.69	7 (22%)
5	ATP	C	601	6	26,33,33	2.23	6 (23%)	31,52,52	1.56	7 (22%)

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
7	ADP	F	501	6	-	5/12/32/32	0/3/3/3
7	ADP	D	501	-	-	4/12/32/32	0/3/3/3
5	ATP	B	601	6	-	3/18/38/38	0/3/3/3
5	ATP	A	601	6	-	2/18/38/38	0/3/3/3
5	ATP	C	601	6	-	5/18/38/38	0/3/3/3

All (26) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	601	ATP	O4'-C1'	7.75	1.51	1.41
5	C	601	ATP	O4'-C1'	7.47	1.51	1.41
5	B	601	ATP	O4'-C1'	7.42	1.51	1.41
7	F	501	ADP	O4'-C1'	6.43	1.50	1.41
5	A	601	ATP	C2'-C1'	-4.56	1.46	1.53
5	C	601	ATP	C2'-C1'	-4.51	1.46	1.53
7	F	501	ADP	C6-N6	4.44	1.50	1.34
5	B	601	ATP	C2'-C1'	-4.27	1.47	1.53
7	F	501	ADP	C2'-C1'	-4.11	1.47	1.53
5	A	601	ATP	C6-N6	3.44	1.46	1.34
7	F	501	ADP	O4'-C4'	3.36	1.52	1.45
5	C	601	ATP	C6-N6	3.36	1.46	1.34
5	B	601	ATP	C6-N6	3.35	1.46	1.34
5	C	601	ATP	C2'-C3'	-3.21	1.44	1.53
5	B	601	ATP	C2'-C3'	-3.18	1.44	1.53
5	A	601	ATP	C2'-C3'	-3.03	1.45	1.53
7	F	501	ADP	C2'-C3'	-2.73	1.45	1.53
7	F	501	ADP	C3'-C4'	-2.62	1.46	1.53
5	B	601	ATP	C4-N3	-2.52	1.32	1.35
5	A	601	ATP	O4'-C4'	2.49	1.50	1.45
5	C	601	ATP	C4-N3	-2.44	1.32	1.35
5	C	601	ATP	O4'-C4'	2.41	1.50	1.45
5	B	601	ATP	O4'-C4'	2.28	1.50	1.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	601	ATP	C4-N3	-2.27	1.32	1.35
7	F	501	ADP	PA-O5'	2.14	1.68	1.59
7	F	501	ADP	C6-C5	-2.10	1.35	1.43

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
5	B	601	ATP	C3'-C2'-C1'	5.08	108.63	100.98
5	A	601	ATP	N3-C2-N1	-4.20	122.11	128.68
7	F	501	ADP	C3'-C2'-C1'	4.19	107.29	100.98
7	F	501	ADP	N3-C2-N1	-4.04	122.36	128.68
5	B	601	ATP	N3-C2-N1	-3.94	122.51	128.68
5	C	601	ATP	N3-C2-N1	-3.93	122.54	128.68
5	A	601	ATP	C4-C5-N7	-3.83	105.41	109.40
5	B	601	ATP	C4-C5-N7	-3.71	105.53	109.40
5	C	601	ATP	C4-C5-N7	-3.64	105.60	109.40
5	A	601	ATP	C3'-C2'-C1'	3.62	106.43	100.98
7	F	501	ADP	C4-C5-N7	-3.61	105.63	109.40
5	A	601	ATP	PB-O3B-PG	-3.24	121.72	132.83
5	B	601	ATP	C2'-C3'-C4'	2.93	108.33	102.64
5	C	601	ATP	C3'-C2'-C1'	2.88	105.32	100.98
7	D	501	ADP	O4'-C1'-C2'	-2.80	102.84	106.93
5	A	601	ATP	C2'-C3'-C4'	2.74	107.97	102.64
5	C	601	ATP	PB-O3B-PG	-2.71	123.53	132.83
5	B	601	ATP	PA-O3A-PB	-2.61	123.88	132.83
7	F	501	ADP	C1'-N9-C4	-2.56	122.14	126.64
5	C	601	ATP	C1'-N9-C4	-2.41	122.41	126.64
5	A	601	ATP	PA-O3A-PB	-2.29	124.96	132.83
5	B	601	ATP	PB-O3B-PG	-2.28	125.02	132.83
5	B	601	ATP	C1'-N9-C4	-2.26	122.67	126.64
5	C	601	ATP	C2'-C3'-C4'	2.23	106.98	102.64
5	A	601	ATP	C1'-N9-C4	-2.21	122.76	126.64
7	D	501	ADP	C5-C6-N6	2.20	123.70	120.35
5	C	601	ATP	PA-O3A-PB	-2.20	125.28	132.83
7	F	501	ADP	PA-O3A-PB	-2.15	125.44	132.83
7	F	501	ADP	C2'-C3'-C4'	2.04	106.60	102.64

There are no chirality outliers.

All (19) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	F	501	ADP	PA-O3A-PB-O3B

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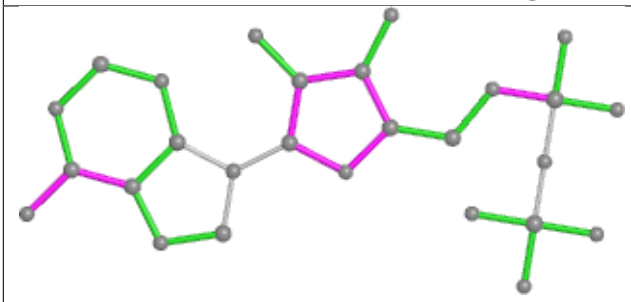
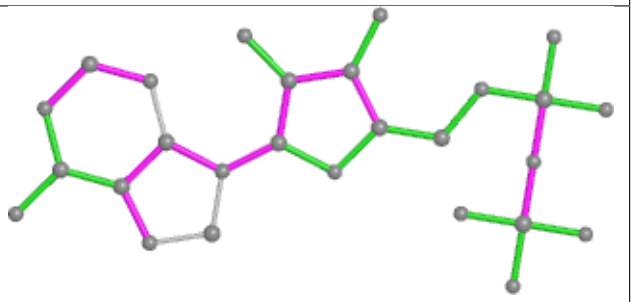
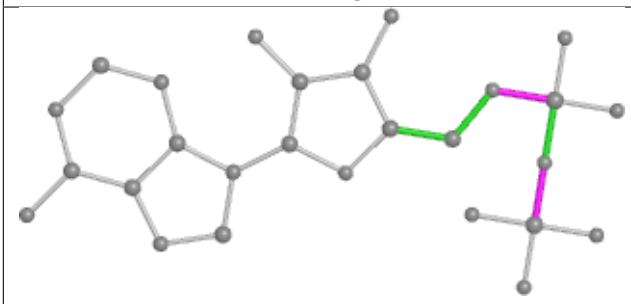
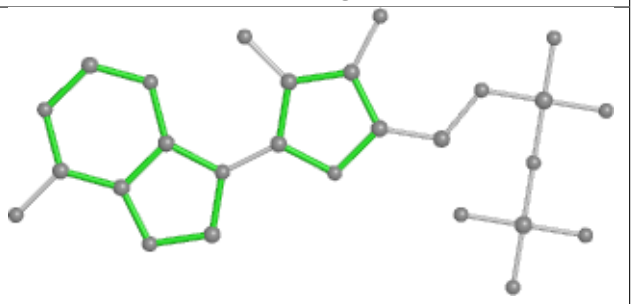
Mol	Chain	Res	Type	Atoms
7	F	501	ADP	C5'-O5'-PA-O3A
7	D	501	ADP	PA-O3A-PB-O2B
7	D	501	ADP	C5'-O5'-PA-O1A
7	D	501	ADP	C5'-O5'-PA-O3A
5	B	601	ATP	O4'-C4'-C5'-O5'
5	C	601	ATP	PB-O3B-PG-O1G
7	F	501	ADP	PA-O3A-PB-O1B
5	A	601	ATP	C5'-O5'-PA-O3A
5	B	601	ATP	PB-O3A-PA-O2A
5	C	601	ATP	C5'-O5'-PA-O1A
5	A	601	ATP	C5'-O5'-PA-O2A
7	F	501	ADP	C5'-O5'-PA-O1A
7	F	501	ADP	C5'-O5'-PA-O2A
7	D	501	ADP	C5'-O5'-PA-O2A
5	C	601	ATP	PB-O3A-PA-O2A
5	C	601	ATP	C5'-O5'-PA-O3A
5	C	601	ATP	PB-O3A-PA-O1A
5	B	601	ATP	C5'-O5'-PA-O1A

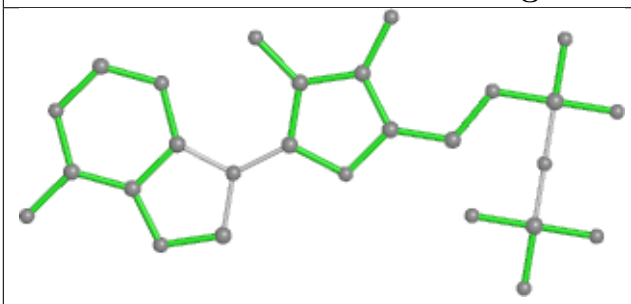
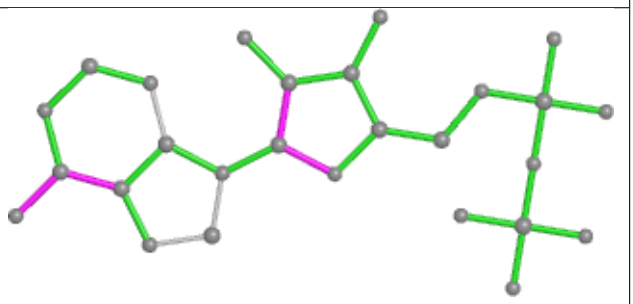
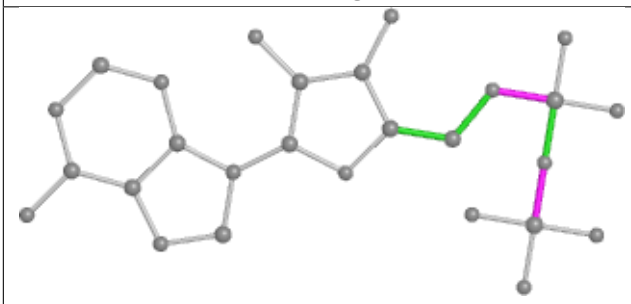
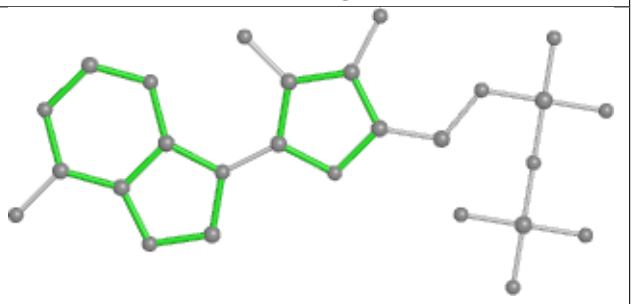
There are no ring outliers.

3 monomers are involved in 9 short contacts:

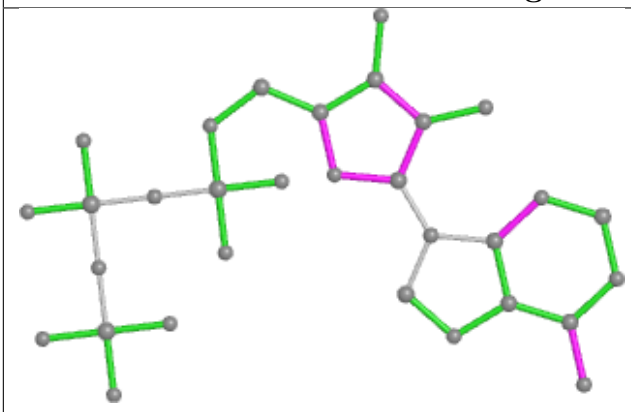
Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	D	501	ADP	6	0
5	B	601	ATP	2	0
5	C	601	ATP	1	0

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.

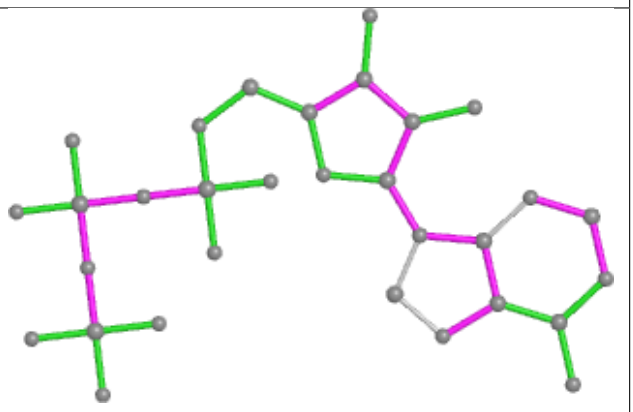
Ligand ADP F 501	
 Bond lengths	 Bond angles
 Torsions	 Rings

Ligand ADP D 501	
 Bond lengths	 Bond angles
 Torsions	 Rings

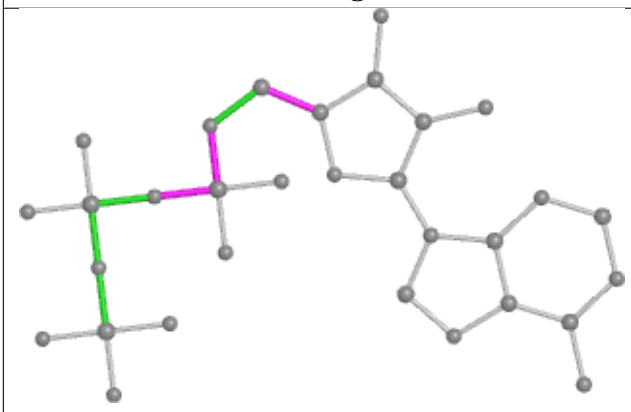
Ligand ATP B 601



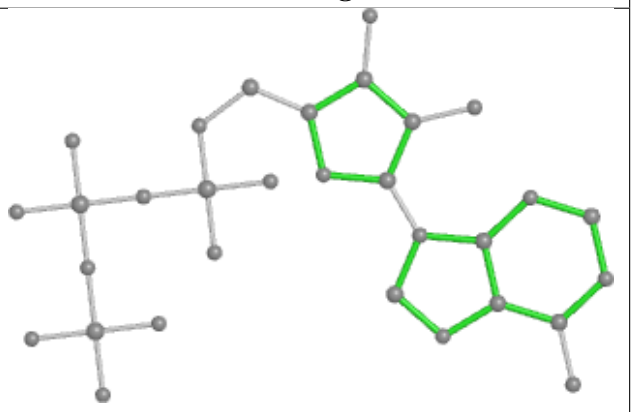
Bond lengths



Bond angles

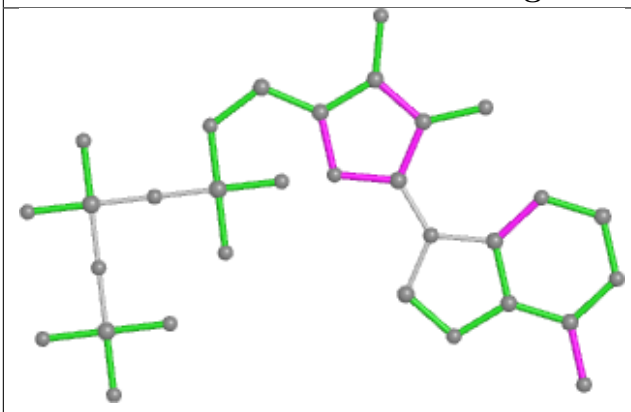


Torsions

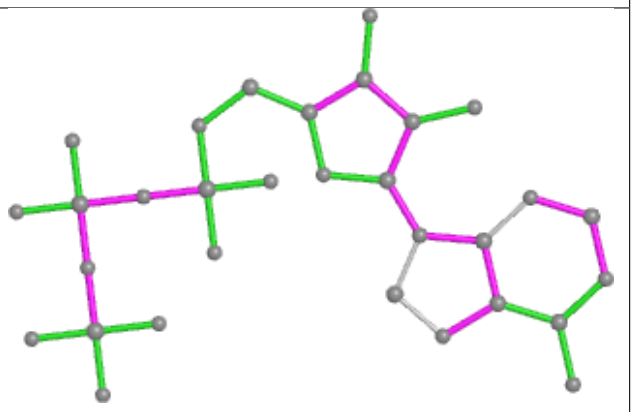


Rings

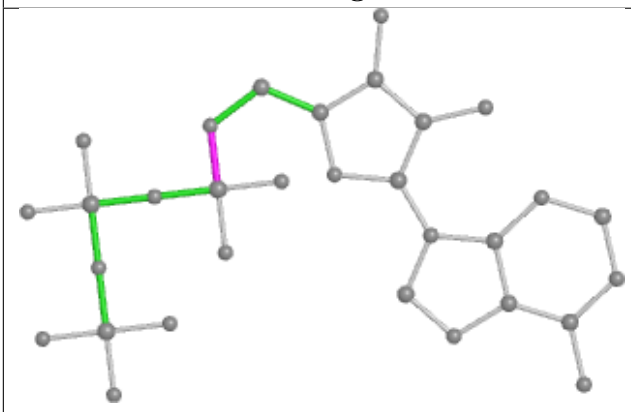
Ligand ATP A 601



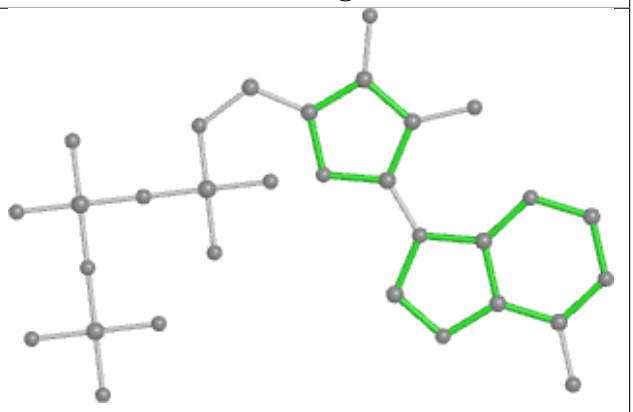
Bond lengths



Bond angles

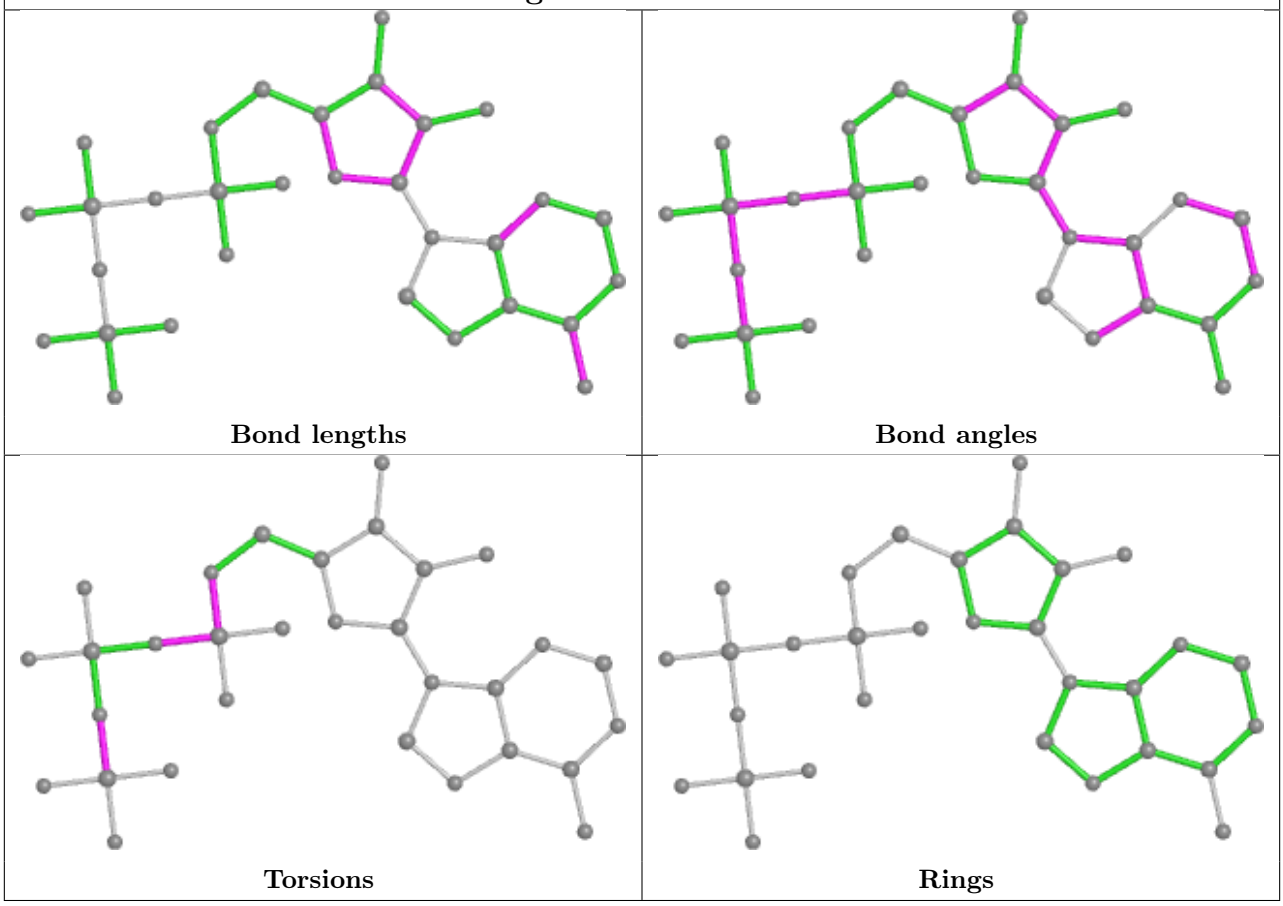


Torsions



Rings

Ligand ATP C 601



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.

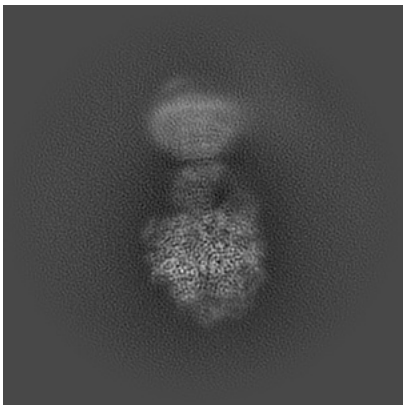
6 Map visualisation

This section contains visualisations of the EMDB entry EMD-34568. These allow visual inspection of the internal detail of the map and identification of artifacts.

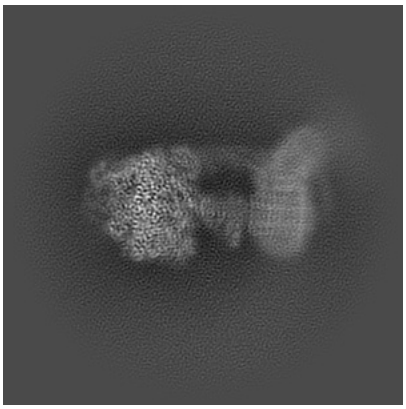
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections

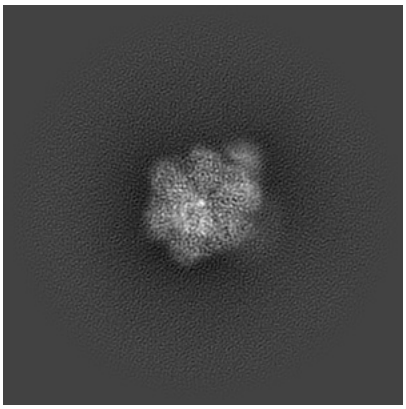
6.1.1 Primary map



X

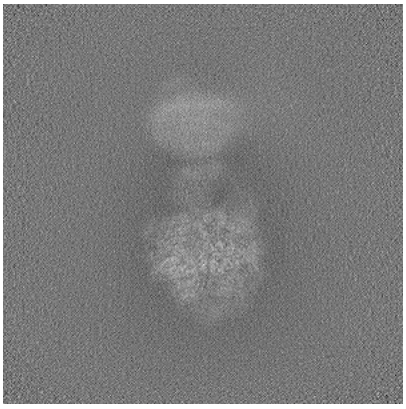


Y

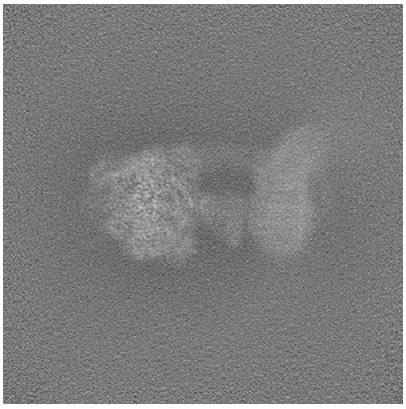


Z

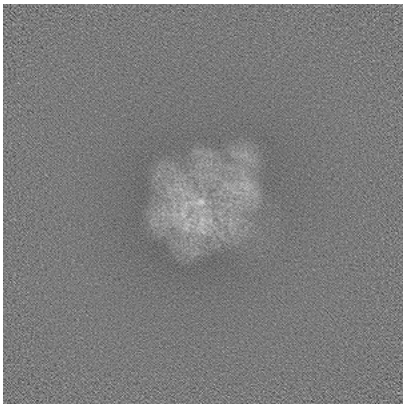
6.1.2 Raw map



X



Y

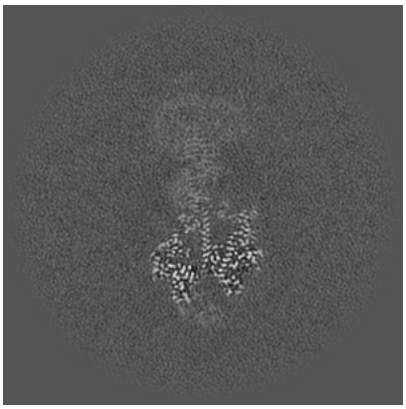


Z

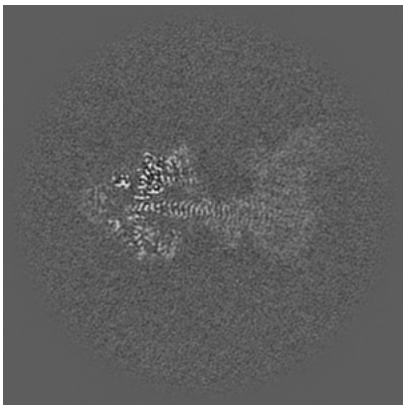
The images above show the map projected in three orthogonal directions.

6.2 Central slices (i)

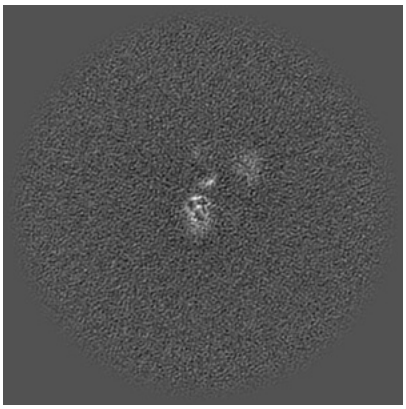
6.2.1 Primary map



X Index: 256

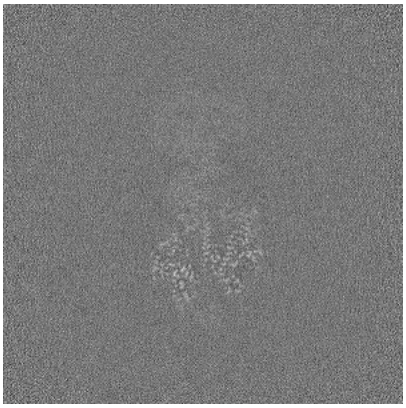


Y Index: 256

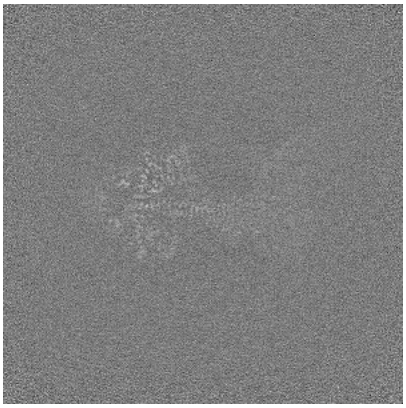


Z Index: 256

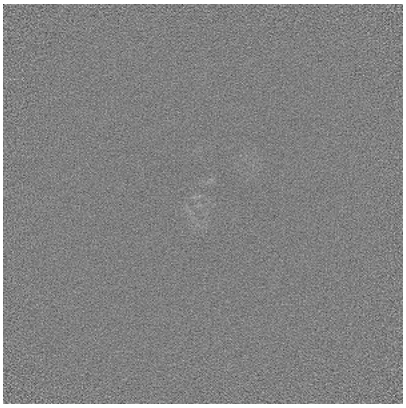
6.2.2 Raw map



X Index: 256



Y Index: 256

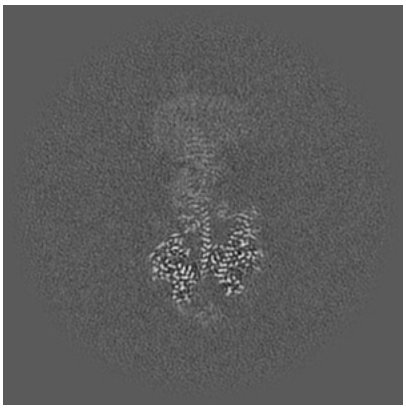


Z Index: 256

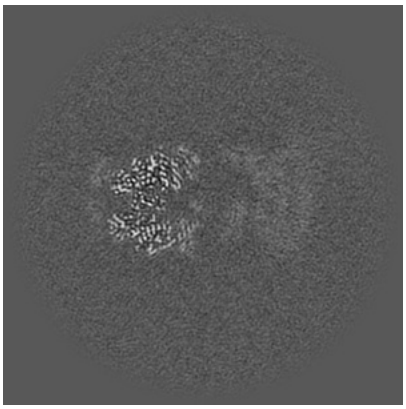
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices (i)

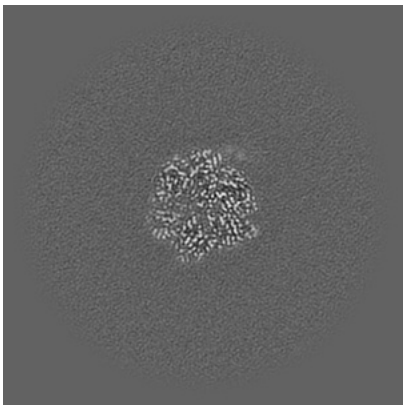
6.3.1 Primary map



X Index: 255

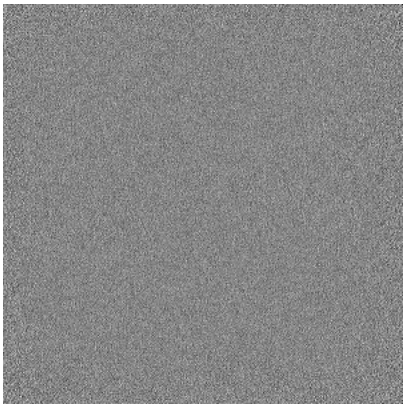


Y Index: 275

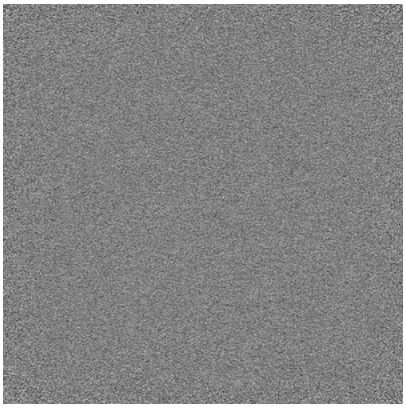


Z Index: 183

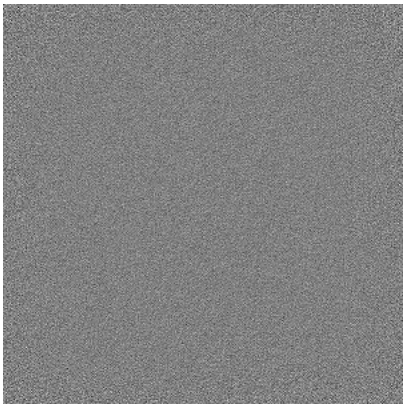
6.3.2 Raw map



X Index: 0



Y Index: 0

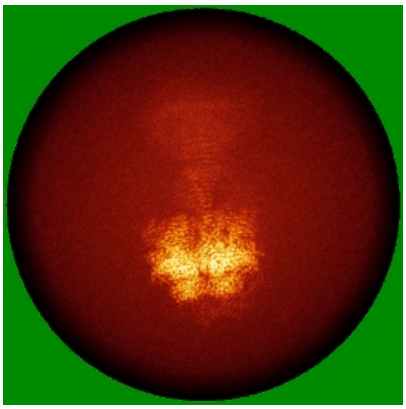


Z Index: 0

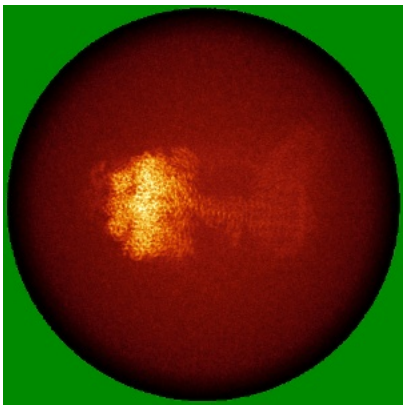
The images above show the largest variance slices of the map in three orthogonal directions.

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

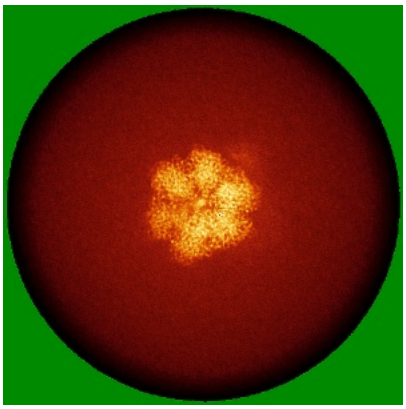
6.4.1 Primary map



X

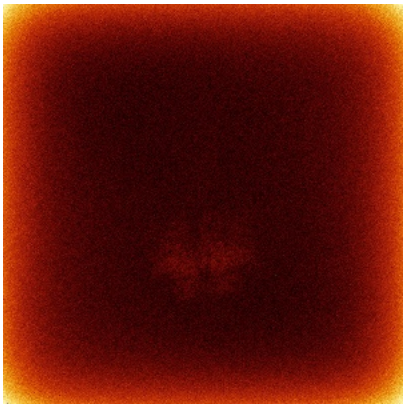


Y

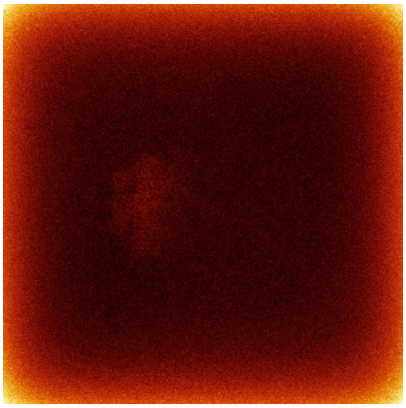


Z

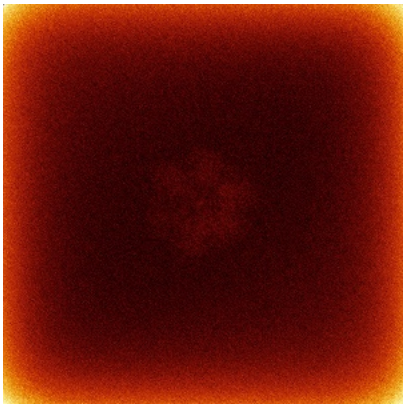
6.4.2 Raw map



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.

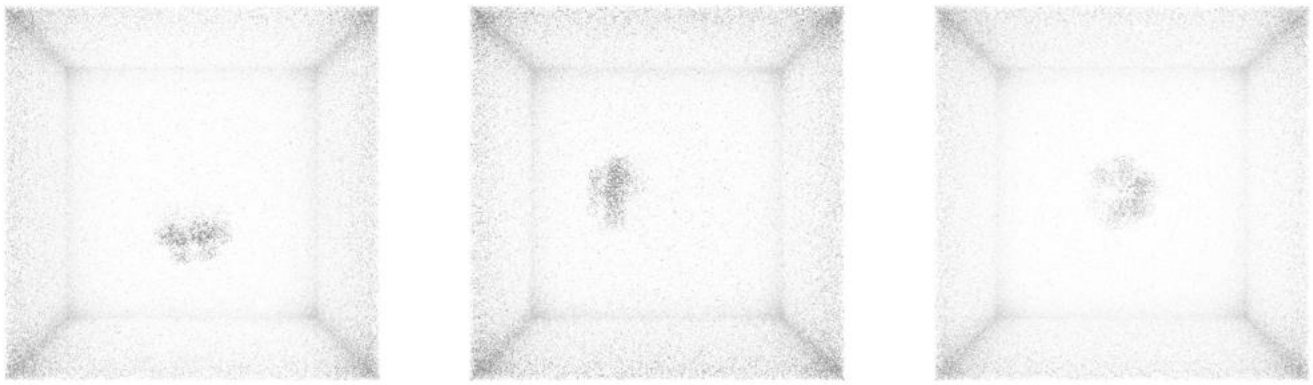
6.5 Orthogonal surface views (i)

6.5.1 Primary map

X Y Z

The images above show the 3D surface view of the map at the recommended contour level 0.1. 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



X Y Z

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.

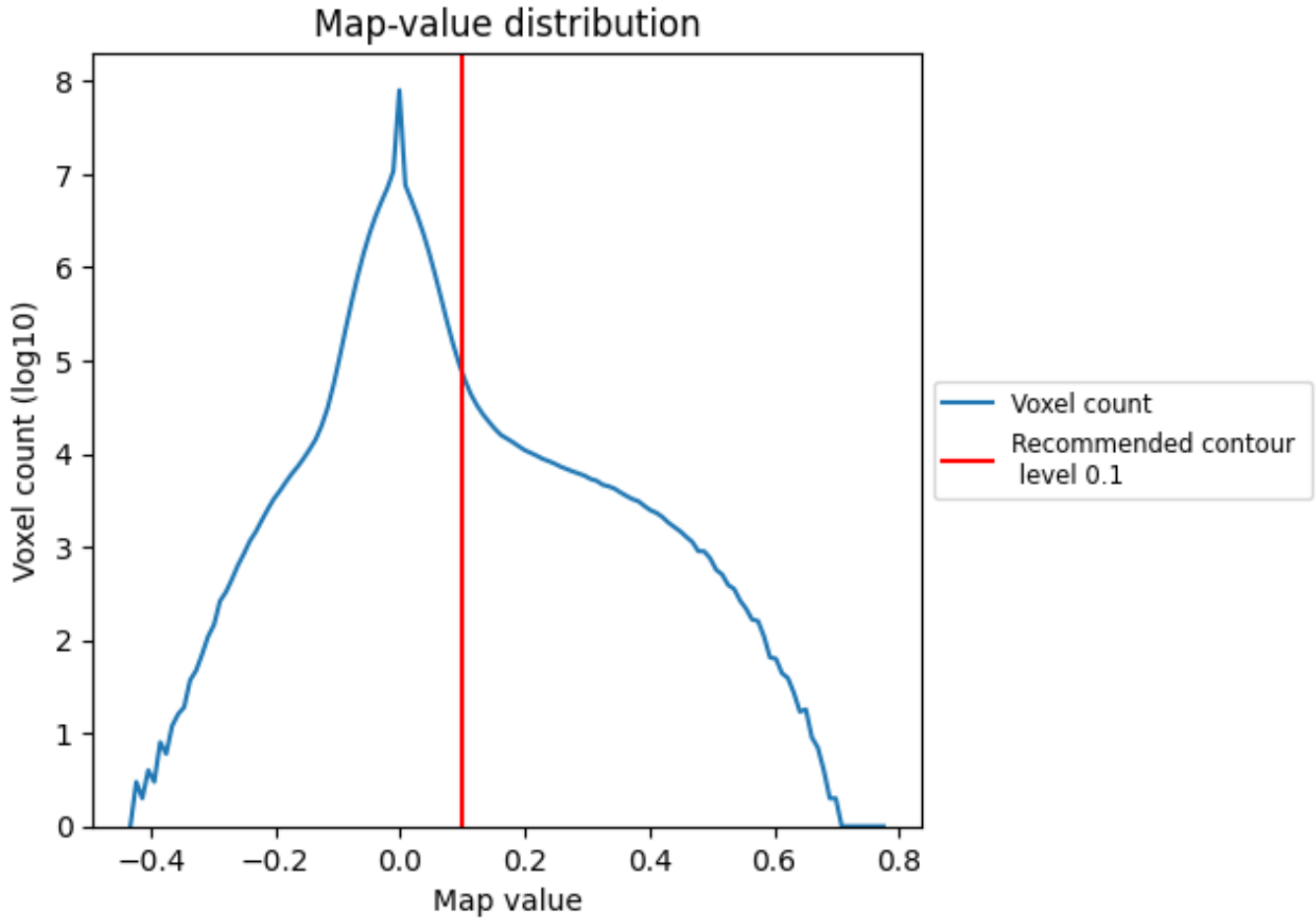
6.6 Mask visualisation (i)

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

7 Map analysis

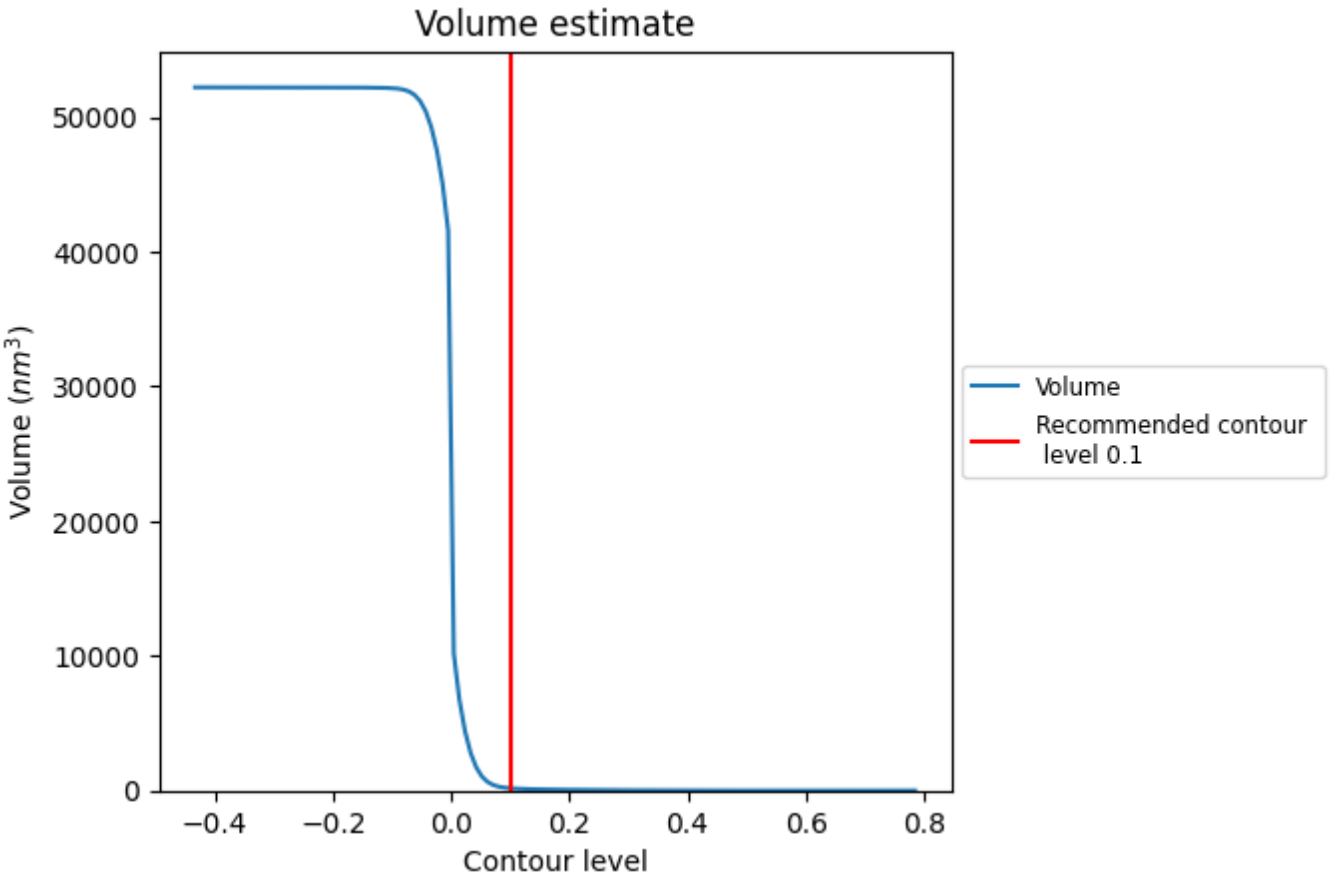
This section contains the results of statistical analysis of the map.

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.

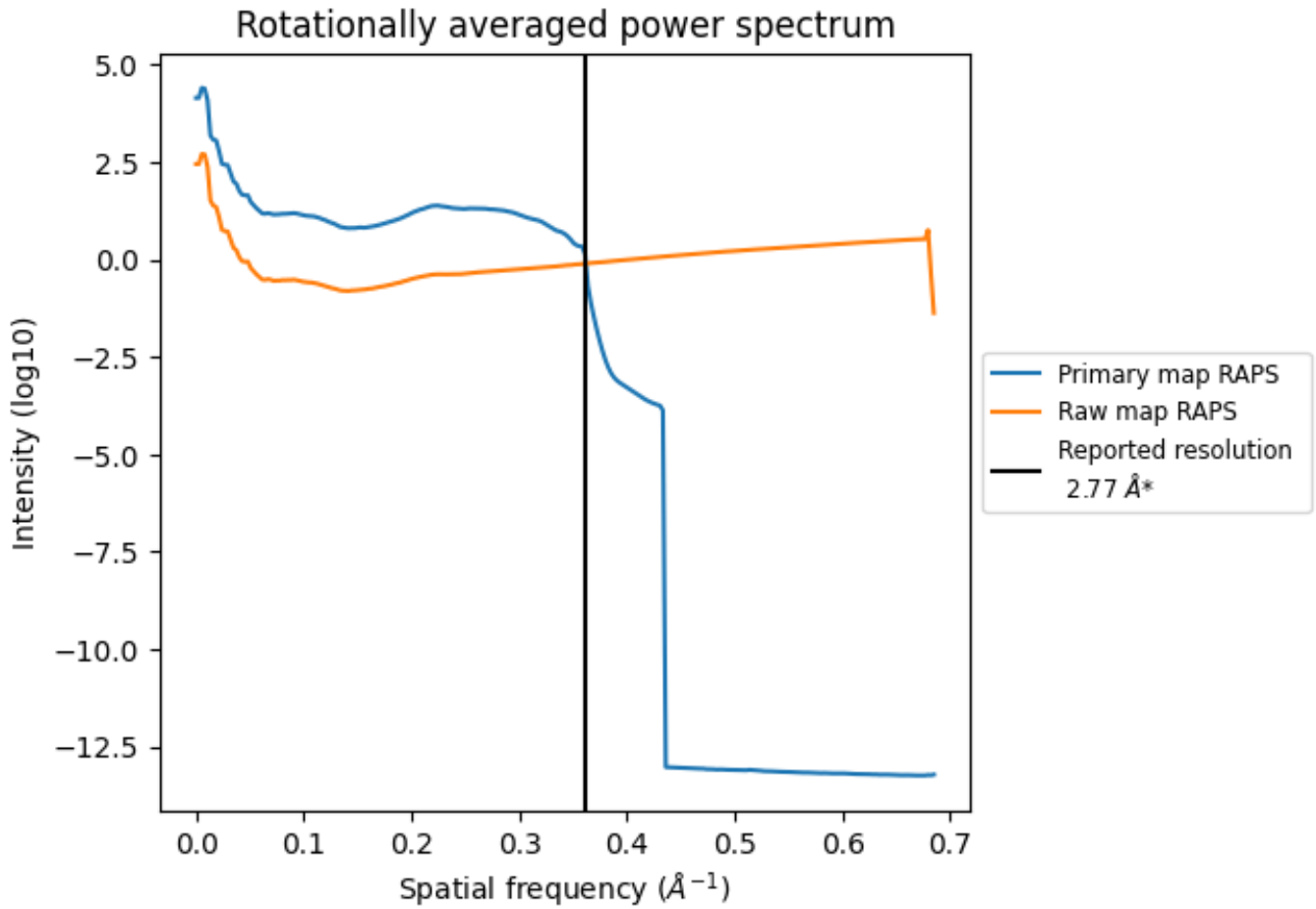
7.2 Volume estimate (i)



The volume at the recommended contour level is 171 nm^3; this corresponds to an approximate mass of 155 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)

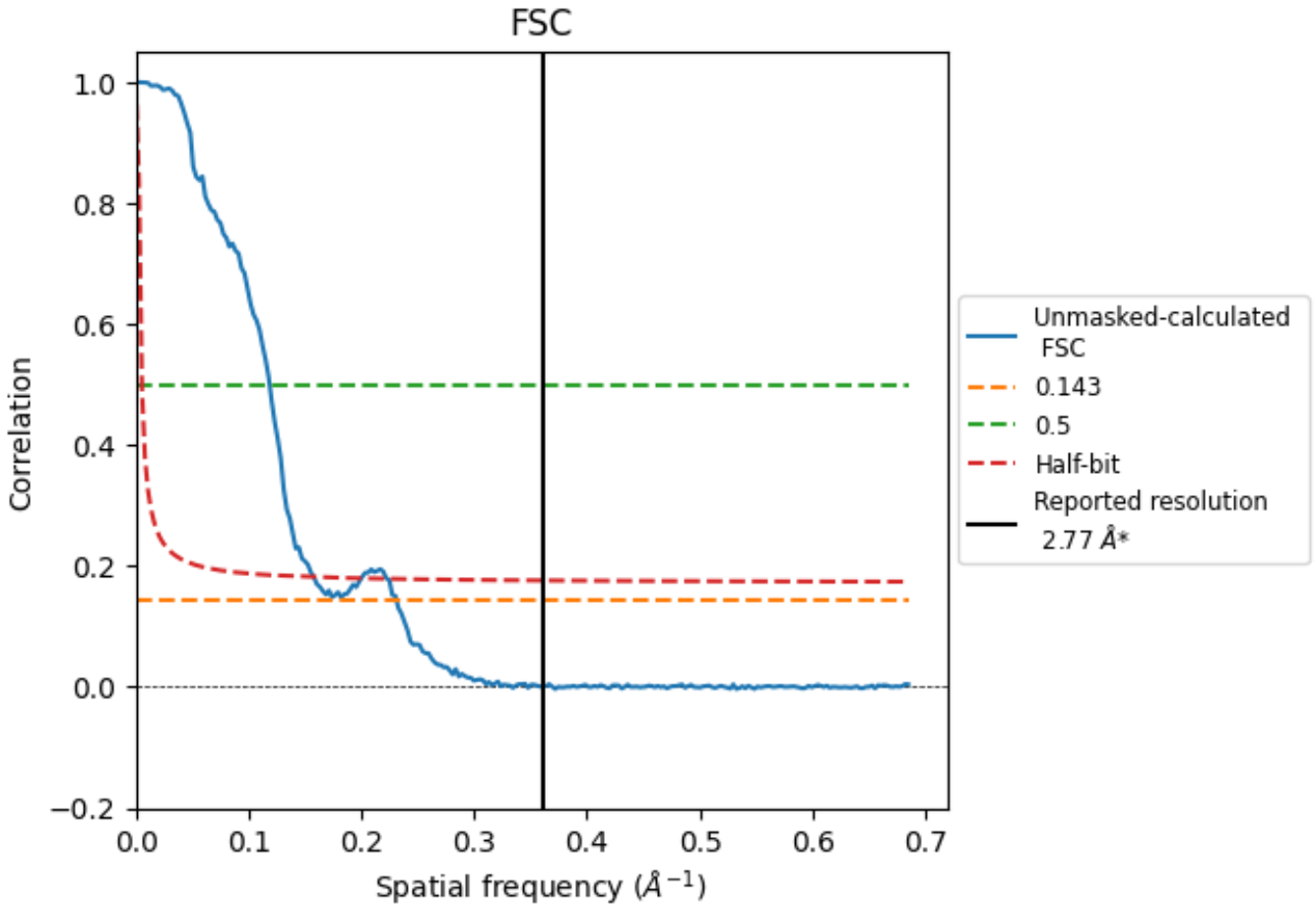


*Reported resolution corresponds to spatial frequency of 0.361 Å^{-1}

8 Fourier-Shell correlation

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



*Reported resolution corresponds to spatial frequency of 0.361 Å^{-1}

8.2 Resolution estimates (i)

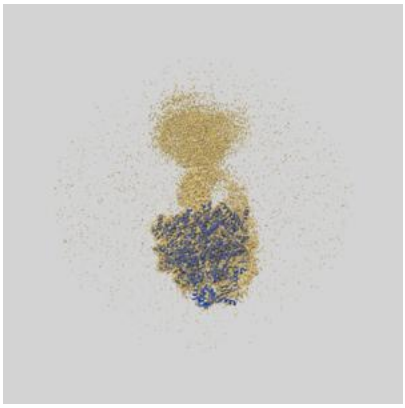
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.77	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	4.33	8.44	6.34

*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 4.33 differs from the reported value 2.77 by more than 10 %

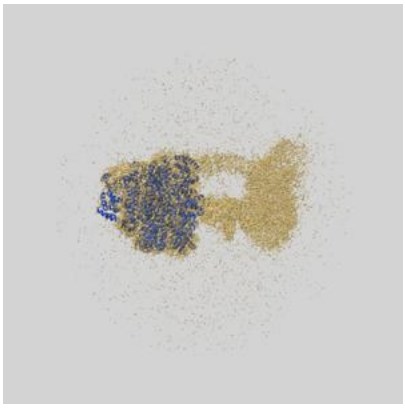
9 Map-model fit

This section contains information regarding the fit between EMDB map EMD-34568 and PDB model 8H9I. Per-residue inclusion information can be found in section 3 on page 6.

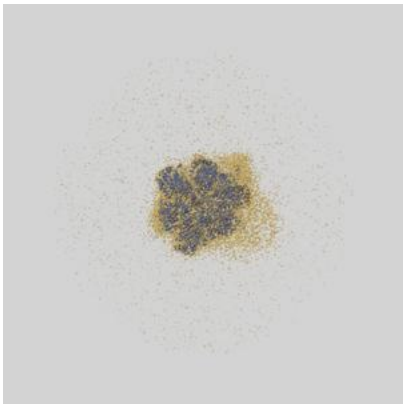
9.1 Map-model overlay



X



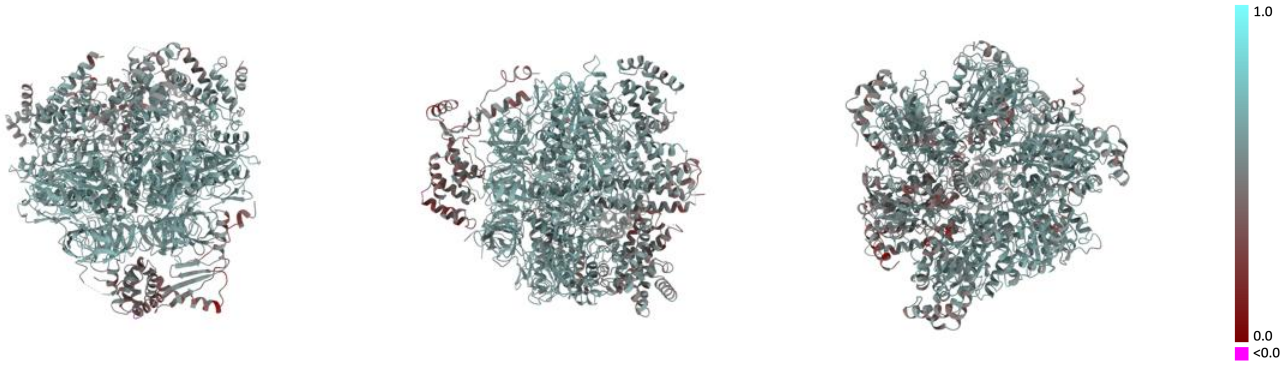
Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.1 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.

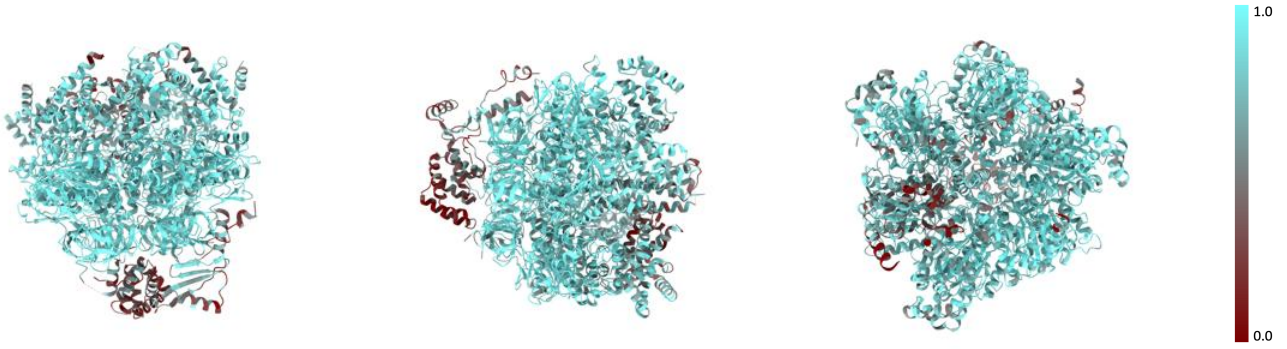
9.2 Q-score mapped to coordinate model (i)



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.

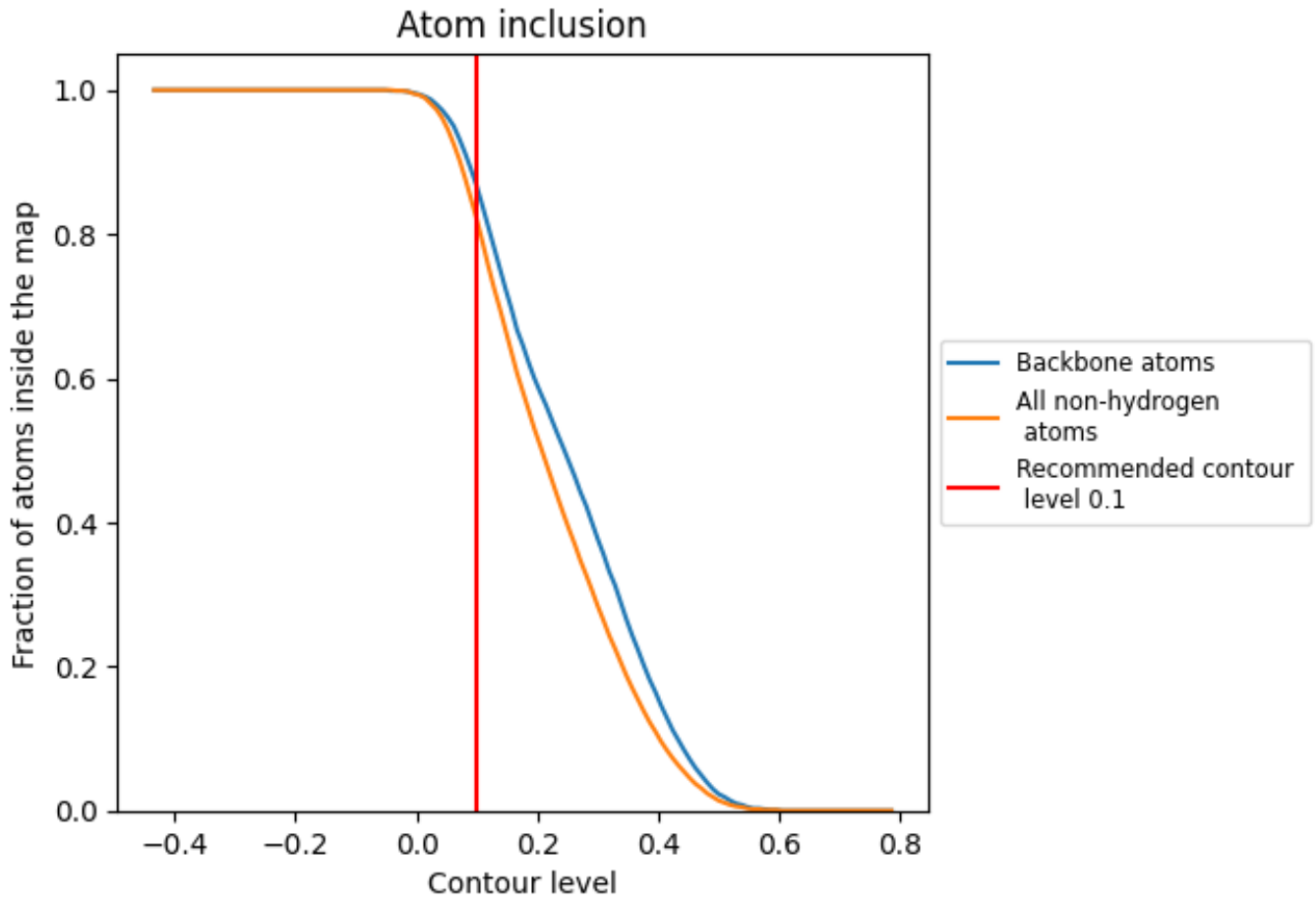
9.3 Atom inclusion mapped to coordinate model (i)



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 (0.1).

9.4 Atom inclusion (i)



At the recommended contour level, 87% of all backbone atoms, 82% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary (i)

The table lists the average atom inclusion at the recommended contour level (0.1) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.8200	0.5640
A	0.8560	0.5710
B	0.8790	0.5890
C	0.8550	0.5740
D	0.7370	0.5300
E	0.8690	0.5800
F	0.9010	0.5970
G	0.7680	0.5460
O	0.3780	0.4160