



Full wwPDB EM Validation Report ⓘ

Mar 12, 2026 – 11:36 PM UTC

PDB ID : 7OIM / pdb_00007oim
EMDB ID : EMD-12932
Title : Mouse RNF213, with mixed nucleotides bound
Authors : Grabarczyk, D.; Ahel, J.; Clausen, T.
Deposited on : 2021-05-11
Resolution : 4.00 Å(reported)
Based on initial model : 6TAX

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

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

A user guide is available at

<https://www.wwpdb.org/validation/2017/EMValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

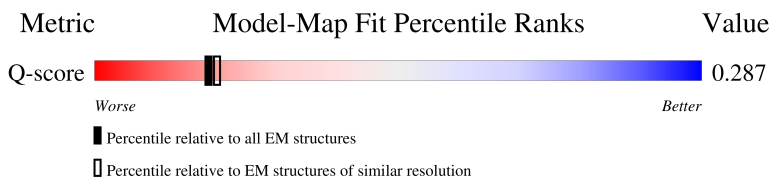
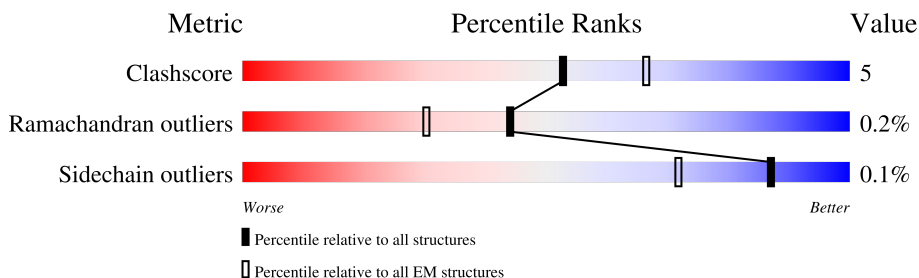
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 4.00 Å.

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



Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	7587 (3.50 - 4.50)

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

Mol	Chain	Length	Quality of chain
1	A	4836	 78% 14% 8%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 35597 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 E3 ubiquitin-protein ligase RNF213.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	4426	Total	C	N	O	S	0	0
			35505	22641	6093	6561	210		

There are 30 discrepancies between the modelled and reference sequences:

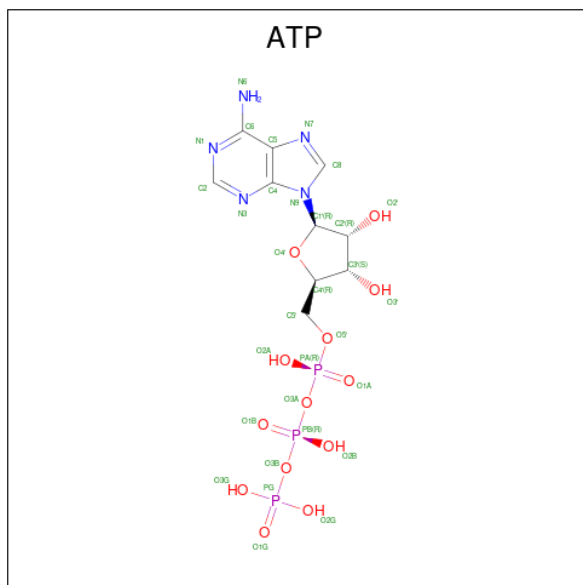
Chain	Residue	Modelled	Actual	Comment	Reference
A	326	MET	-	initiating methionine	UNP E9Q555
A	327	ALA	-	expression tag	UNP E9Q555
A	328	SER	-	expression tag	UNP E9Q555
A	329	TRP	-	expression tag	UNP E9Q555
A	330	SER	-	expression tag	UNP E9Q555
A	331	HIS	-	expression tag	UNP E9Q555
A	332	PRO	-	expression tag	UNP E9Q555
A	333	GLN	-	expression tag	UNP E9Q555
A	334	PHE	-	expression tag	UNP E9Q555
A	335	GLU	-	expression tag	UNP E9Q555
A	336	LYS	-	expression tag	UNP E9Q555
A	337	GLY	-	expression tag	UNP E9Q555
A	338	SER	-	expression tag	UNP E9Q555
A	?	-	VAL	deletion	UNP E9Q555
A	?	-	ARG	deletion	UNP E9Q555
A	?	-	ASN	deletion	UNP E9Q555
A	?	-	ARG	deletion	UNP E9Q555
A	5149	GLY	-	expression tag	UNP E9Q555
A	5150	GLY	-	expression tag	UNP E9Q555
A	5151	GLY	-	expression tag	UNP E9Q555
A	5152	HIS	-	expression tag	UNP E9Q555
A	5153	HIS	-	expression tag	UNP E9Q555
A	5154	HIS	-	expression tag	UNP E9Q555
A	5155	HIS	-	expression tag	UNP E9Q555
A	5156	HIS	-	expression tag	UNP E9Q555
A	5157	HIS	-	expression tag	UNP E9Q555
A	5158	HIS	-	expression tag	UNP E9Q555
A	5159	HIS	-	expression tag	UNP E9Q555

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Chain	Residue	Modelled	Actual	Comment	Reference
A	5160	HIS	-	expression tag	UNP E9Q555
A	5161	HIS	-	expression tag	UNP E9Q555

- Molecule 2 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
2	A	1	Total	C	N	O	P	0
			31	10	5	13	3	

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		AltConf
3	A	1	Total	Mg	0
			1	1	

- Molecule 4 is ZINC ION (CCD ID: ZN) (formula: Zn) (labeled as "Ligand of Interest" by depositor).

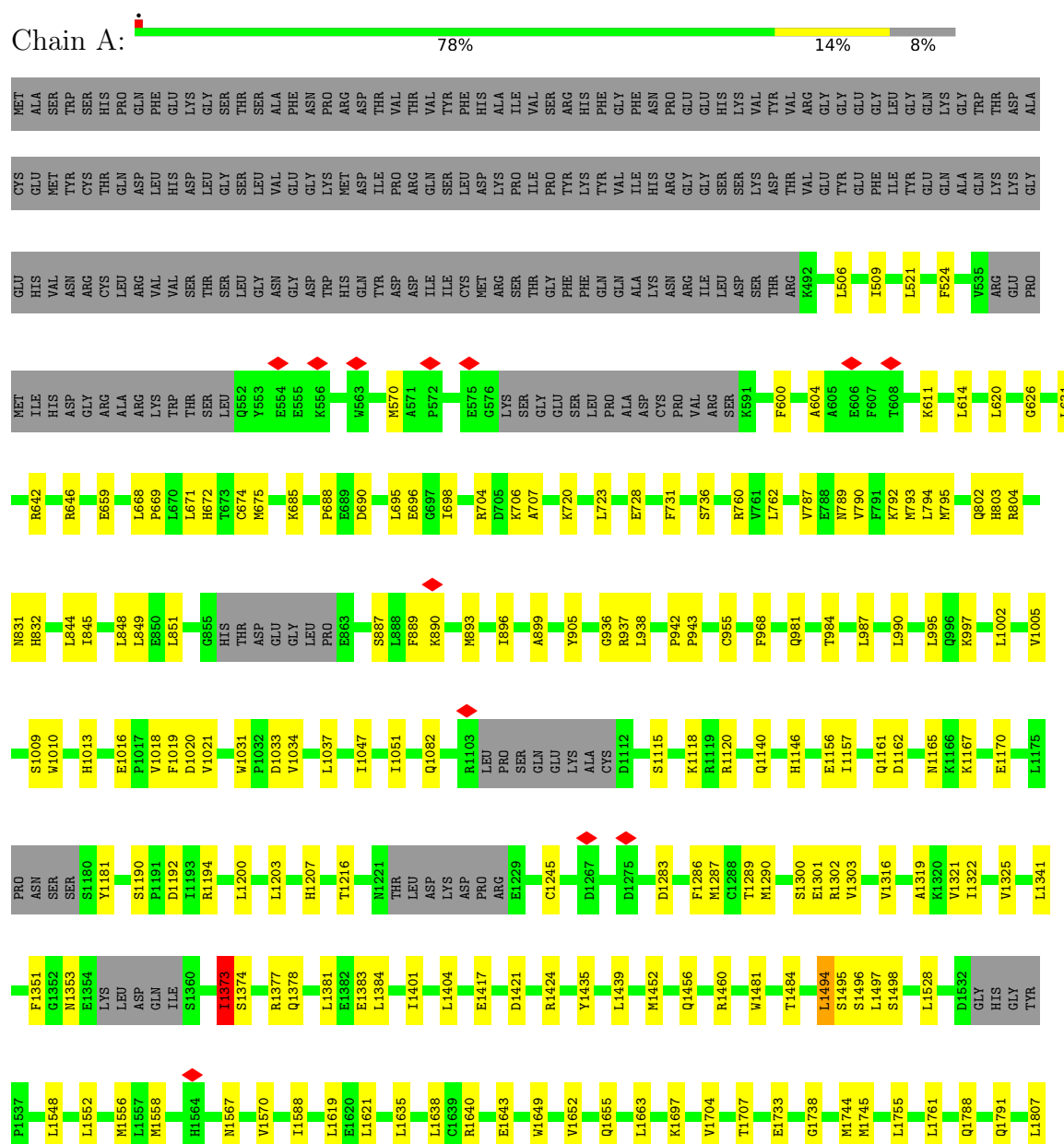
Mol	Chain	Residues	Atoms		AltConf
4	A	2	Total	Zn	0
			2	2	

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

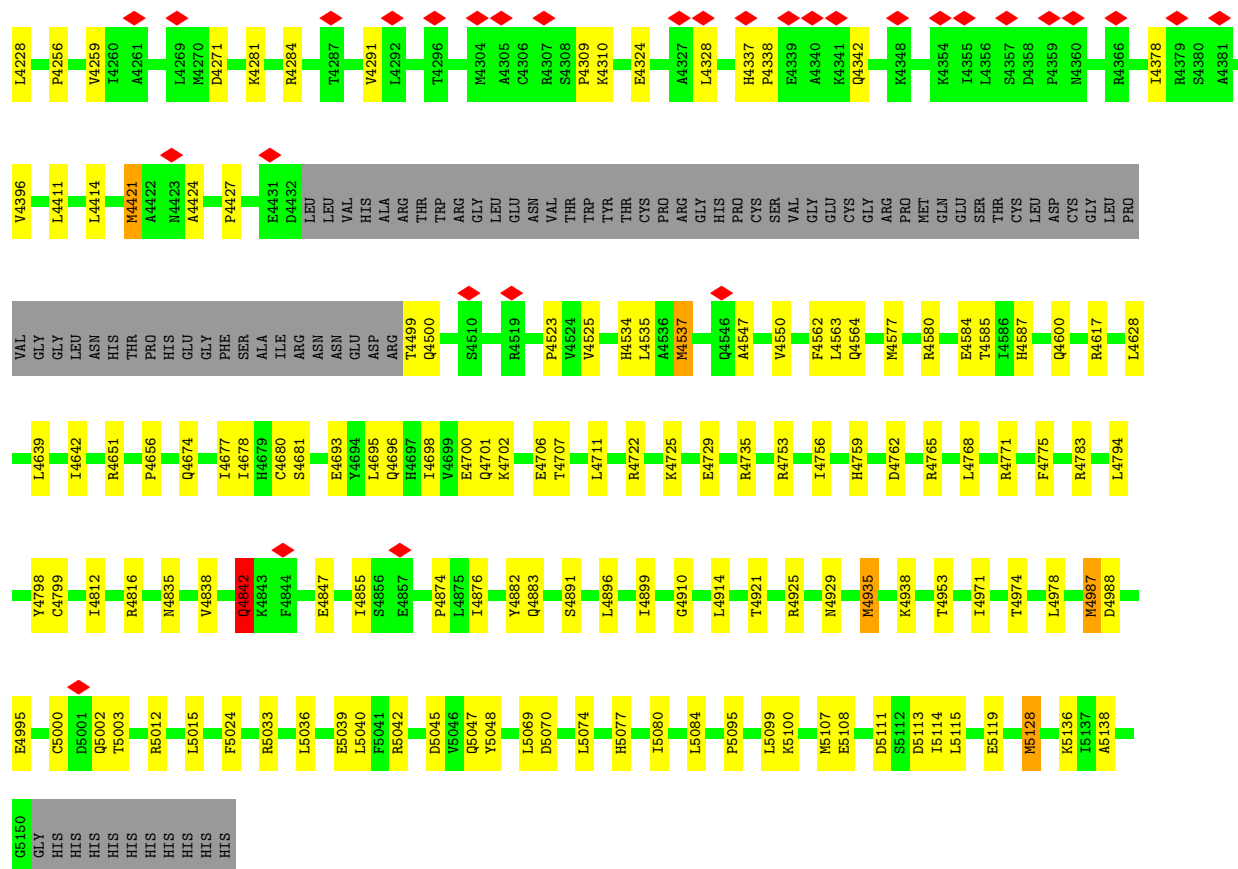
3 Residue-property plots

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: E3 ubiquitin-protein ligase RNF213







4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	336000	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	TFS KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	45, 40.5	Depositor
Minimum defocus (nm)	-1500	Depositor
Maximum defocus (nm)	-2500	Depositor
Magnification	Not provided	
Image detector	GATAN K3 BIOQUANTUM (6k x 4k), GATAN K2 QUANTUM (4k x 4k)	Depositor
Maximum map value	0.034	Depositor
Minimum map value	-0.010	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.001	Depositor
Recommended contour level	0.005	Depositor
Map size (Å)	370.656, 370.656, 370.656	wwPDB
Map dimensions	352, 352, 352	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.053, 1.053, 1.053	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ATP, ZN, MG, AGS, 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	0.27	0/36242	0.72	49/49026 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	6

There are no bond length outliers.

All (49) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1494	LEU	O-C-N	10.89	137.07	122.59
1	A	4701	GLN	CA-CB-CG	8.97	132.05	114.10
1	A	3121	PRO	CA-C-N	7.73	136.42	122.13
1	A	3121	PRO	C-N-CA	7.73	136.42	122.13
1	A	4701	GLN	N-CA-CB	7.40	121.11	110.16
1	A	4421	MET	CA-CB-CG	7.30	128.69	114.10
1	A	5108	GLU	N-CA-CB	6.99	121.17	110.61
1	A	3864	GLN	N-CA-CB	6.90	120.26	110.12
1	A	1383	GLU	N-CA-CB	6.71	120.68	110.28
1	A	3415	THR	O-C-N	-6.40	114.86	122.15
1	A	5108	GLU	CA-CB-CG	6.22	126.55	114.10
1	A	1013	HIS	CA-C-N	6.15	133.40	123.93
1	A	1013	HIS	C-N-CA	6.15	133.40	123.93
1	A	4309	PRO	CA-C-N	6.08	136.64	121.80
1	A	4309	PRO	C-N-CA	6.08	136.64	121.80
1	A	671	LEU	CA-CB-CG	5.89	136.93	116.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	4987	MET	CA-CB-CG	5.88	125.86	114.10
1	A	4700	GLU	CB-CG-CD	5.81	122.48	112.60
1	A	2369	MET	CB-CG-SD	5.73	129.90	112.70
1	A	4002	ARG	CB-CG-CD	5.71	124.43	111.30
1	A	762	LEU	CA-CB-CG	5.66	136.12	116.30
1	A	2878	ILE	N-CA-C	-5.66	106.87	111.91
1	A	3909	LEU	N-CA-C	5.66	120.18	113.28
1	A	4842	GLN	CA-CB-CG	5.65	125.41	114.10
1	A	5108	GLU	CB-CA-C	-5.63	100.65	110.17
1	A	3910	GLU	CA-CB-CG	5.58	125.25	114.10
1	A	3495	GLN	CA-C-N	5.52	132.08	121.54
1	A	3495	GLN	C-N-CA	5.52	132.08	121.54
1	A	4002	ARG	N-CA-CB	5.50	118.05	110.07
1	A	4310	LYS	N-CA-C	5.48	121.92	109.81
1	A	4537	MET	CA-CB-CG	5.47	125.05	114.10
1	A	1975	LEU	CA-CB-CG	5.43	135.29	116.30
1	A	2025	GLN	N-CA-C	5.42	118.69	111.75
1	A	4995	GLU	CA-CB-CG	5.40	124.89	114.10
1	A	1934	ARG	CA-CB-CG	5.38	124.86	114.10
1	A	3135	MET	CB-CG-SD	-5.33	96.70	112.70
1	A	1377	ARG	CA-CB-CG	5.32	124.73	114.10
1	A	4847	GLU	CA-C-N	5.32	131.69	121.54
1	A	4847	GLU	C-N-CA	5.32	131.69	121.54
1	A	4700	GLU	N-CA-CB	5.31	118.50	110.28
1	A	4935	MET	CB-CG-SD	5.30	128.61	112.70
1	A	1051	ILE	CA-C-N	5.17	131.42	121.54
1	A	1051	ILE	C-N-CA	5.17	131.42	121.54
1	A	4700	GLU	CA-C-N	-5.15	112.98	120.29
1	A	4700	GLU	C-N-CA	-5.15	112.98	120.29
1	A	3865	GLU	N-CA-CB	5.14	118.25	110.28
1	A	1373	ILE	CA-C-N	5.07	131.23	121.54
1	A	1373	ILE	C-N-CA	5.07	131.23	121.54
1	A	1452	MET	CB-CG-SD	5.07	127.91	112.70

There are no chirality outliers.

All (6) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	1016	GLU	Peptide
1	A	1170	GLU	Peptide
1	A	1373	ILE	Peptide
1	A	4165	GLU	Peptide

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Mol	Chain	Res	Type	Group
1	A	659	GLU	Peptide
1	A	688	PRO	Peptide

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	35505	0	35642	378	0
2	A	31	0	12	3	0
3	A	1	0	0	0	0
4	A	2	0	0	0	0
5	A	27	0	12	0	0
6	A	31	0	12	2	0
All	All	35597	0	35678	378	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 5.

All (378) 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:A:731:PHE:HB3	1:A:760:ARG:HH21	1.47	0.78
1:A:4935:MET:SD	1:A:4938:LYS:NZ	2.60	0.74
1:A:942:PRO:HB3	1:A:3207:ARG:HE	1.54	0.72
1:A:3292:LEU:O	1:A:3296:LEU:HB2	1.91	0.70
1:A:942:PRO:HB3	1:A:3207:ARG:HH21	1.56	0.70
1:A:1216:THR:HG21	1:A:1245:CYS:HB2	1.77	0.66
1:A:4695:LEU:HD11	1:A:4876:ILE:HG12	1.78	0.66
1:A:646:ARG:NH2	1:A:674:CYS:SG	2.69	0.65
1:A:1301:GLU:OE2	1:A:1302:ARG:NH1	2.29	0.65
1:A:2186:GLN:NE2	1:A:2369:MET:SD	2.70	0.65
1:A:4427:PRO:HA	1:A:4617:ARG:HE	1.61	0.64
1:A:728:GLU:O	1:A:760:ARG:NH2	2.30	0.64
1:A:981:GLN:NE2	1:A:3206:TYR:OH	2.30	0.64
1:A:942:PRO:HB3	1:A:3207:ARG:NH2	2.11	0.64
1:A:4185:LEU:HD22	1:A:4228:LEU:HD12	1.80	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:5040:LEU:HD21	1:A:5136:LYS:HE2	1.81	0.63
1:A:4009:ALA:O	1:A:4013:HIS:ND1	2.28	0.63
1:A:4842:GLN:HE21	1:A:4842:GLN:HA	1.65	0.62
1:A:3605:ALA:O	1:A:3609:ASP:HB2	1.99	0.62
1:A:4525:VAL:HG13	1:A:4628:LEU:HD12	1.82	0.62
1:A:2594:CYS:SG	1:A:2595:GLY:N	2.73	0.61
1:A:3780:ARG:NH2	1:A:4910:GLY:O	2.33	0.61
1:A:794:LEU:HD22	1:A:844:LEU:HD11	1.81	0.61
1:A:3064:LEU:HD22	1:A:3109:VAL:HG11	1.83	0.61
1:A:1417:GLU:OE1	1:A:1496:SER:OG	2.07	0.61
1:A:3658:SER:HB3	1:A:3720:LEU:HD11	1.82	0.61
1:A:3091:GLY:HA3	1:A:3096:ARG:HD3	1.83	0.61
1:A:3897:LEU:HB2	1:A:3938:LYS:HE3	1.83	0.61
1:A:4200:GLU:HA	1:A:4206:ARG:HH21	1.65	0.61
1:A:942:PRO:HB3	1:A:3207:ARG:NE	2.15	0.61
1:A:3937:LYS:HG3	1:A:3942:PHE:HA	1.83	0.61
1:A:887:SER:O	1:A:890:LYS:NZ	2.34	0.60
1:A:2294:ILE:HD11	1:A:2299:GLY:HA2	1.82	0.60
1:A:1744:MET:HG3	1:A:1745:MET:HE2	1.85	0.59
1:A:2334:LYS:HD3	1:A:2356:LEU:HD22	1.84	0.59
1:A:893:MET:HE1	1:A:905:TYR:HD2	1.67	0.59
1:A:4735:ARG:NH2	1:A:4855:ILE:O	2.36	0.59
1:A:2128:PRO:HB3	1:A:2183:LEU:HD23	1.84	0.59
1:A:3530:ARG:HH12	1:A:3618:ILE:HA	1.67	0.59
1:A:1384:LEU:HD11	1:A:1439:LEU:HD11	1.84	0.58
1:A:3882:HIS:NE2	1:A:4021:ASP:OD2	2.36	0.58
1:A:3006:THR:HG21	1:A:3011:ALA:HB2	1.86	0.58
1:A:3501:ARG:HH12	1:A:3593:GLY:HA2	1.68	0.58
1:A:942:PRO:CB	1:A:3207:ARG:HH21	2.16	0.58
1:A:4411:LEU:HD23	1:A:4414:LEU:HD12	1.86	0.58
1:A:2459:ILE:HA	1:A:2462:ILE:HD12	1.85	0.58
1:A:893:MET:SD	1:A:937:ARG:NH2	2.78	0.57
1:A:4221:ASN:O	1:A:4224:HIS:ND1	2.36	0.57
1:A:4971:ILE:HA	1:A:4974:THR:HG22	1.85	0.57
1:A:704:ARG:NH2	1:A:736:SER:O	2.37	0.57
1:A:845:ILE:HA	1:A:848:LEU:HG	1.87	0.57
1:A:611:LYS:HA	1:A:614:LEU:HD13	1.87	0.56
1:A:987:LEU:HA	1:A:990:LEU:HD13	1.86	0.56
1:A:2100:ILE:HD13	1:A:2159:LEU:HD23	1.86	0.56
1:A:4759:HIS:O	1:A:4765:ARG:NH1	2.33	0.56
1:A:4142:ASN:ND2	1:A:4144:GLU:OE2	2.38	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:4768:LEU:HD22	1:A:4771:ARG:HH21	1.70	0.56
1:A:2601:ASP:O	1:A:2605:CYS:HB2	2.05	0.56
1:A:4378:ILE:O	1:A:4564:GLN:NE2	2.38	0.56
1:A:2474:HIS:HB2	1:A:2477:SER:HB3	1.87	0.55
1:A:3696:VAL:O	1:A:3700:GLN:NE2	2.30	0.55
1:A:3382:MET:HE3	1:A:5036:LEU:CD1	2.37	0.55
1:A:2590:ARG:NH1	1:A:2660:GLU:OE2	2.40	0.55
1:A:3127:ARG:NH2	6:A:5206:AGS:S1G	2.80	0.55
1:A:936:GLY:O	1:A:3238:THR:OG1	2.25	0.54
1:A:2021:SER:OG	1:A:2022:THR:N	2.40	0.54
1:A:1845:THR:O	1:A:1883:ARG:NH1	2.39	0.54
1:A:4117:LYS:HE3	1:A:4677:ILE:HG12	1.90	0.54
1:A:831:ASN:O	1:A:832:HIS:ND1	2.40	0.54
1:A:1435:TYR:HB3	1:A:1439:LEU:HD13	1.90	0.54
1:A:3479:ARG:NH2	1:A:5000:CYS:SG	2.76	0.54
1:A:5077:HIS:HA	1:A:5080:ILE:HG12	1.89	0.54
1:A:3891:ILE:HG21	1:A:4007:LYS:HG2	1.89	0.54
1:A:3728:GLU:HA	1:A:3731:MET:HG3	1.89	0.53
1:A:4164:GLN:HG2	1:A:4171:SER:H	1.73	0.53
1:A:4835:ASN:ND2	1:A:4914:LEU:O	2.41	0.53
1:A:1162:ASP:O	1:A:1165:ASN:ND2	2.42	0.53
1:A:803:HIS:HB3	1:A:851:LEU:HG	1.91	0.53
1:A:1494:LEU:HG	1:A:1498:SER:OG	2.09	0.53
1:A:2222:ASP:HA	1:A:2258:ARG:HH22	1.74	0.53
1:A:695:LEU:HD22	1:A:698:ILE:HD12	1.91	0.53
1:A:1321:VAL:HG11	1:A:1378:GLN:HG3	1.90	0.53
1:A:4729:GLU:HG3	1:A:4771:ARG:HG2	1.91	0.53
1:A:1181:TYR:OH	1:A:1194:ARG:NH2	2.43	0.52
1:A:3281:SER:OG	1:A:3282:ARG:N	2.42	0.52
1:A:2370:ARG:HB3	1:A:2377:VAL:HG22	1.91	0.52
1:A:2708:ALA:HA	1:A:2863:ARG:HH22	1.74	0.52
1:A:4378:ILE:HG13	1:A:4564:GLN:HE21	1.74	0.52
1:A:2733:GLY:HA3	1:A:2914:LEU:HB2	1.92	0.52
1:A:984:THR:HG21	1:A:3203:GLU:HG3	1.90	0.52
1:A:2162:PHE:HB3	1:A:2176:LEU:HD11	1.90	0.52
1:A:2306:VAL:HG23	1:A:2307:MET:HG3	1.91	0.52
1:A:4783:ARG:NH1	1:A:4799:CYS:O	2.42	0.52
1:A:4842:GLN:HE21	1:A:4842:GLN:CA	2.22	0.52
1:A:1316:VAL:HA	1:A:1341:LEU:HD21	1.92	0.51
1:A:3017:GLN:NE2	1:A:3394:GLY:O	2.42	0.51
1:A:4978:LEU:HD22	1:A:5015:LEU:HD22	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:4175:LEU:HA	1:A:4178:VAL:HG22	1.90	0.51
1:A:2225:THR:O	1:A:2258:ARG:NH1	2.43	0.51
1:A:2142:ASP:OD1	1:A:2142:ASP:N	2.42	0.51
1:A:3960:LEU:HB2	1:A:3964:HIS:HB2	1.92	0.51
1:A:4988:ASP:N	1:A:4988:ASP:OD1	2.43	0.51
1:A:3224:ASP:OD2	1:A:3381:ARG:NH1	2.43	0.51
1:A:4328:LEU:HD13	1:A:4337:HIS:HB3	1.91	0.51
1:A:1156:GLU:HG2	1:A:1157:ILE:HG12	1.92	0.51
1:A:3732:PHE:HA	1:A:3735:MET:HE3	1.92	0.51
1:A:3403:ASP:OD2	1:A:3567:ARG:HD3	2.11	0.51
1:A:2042:MET:HG3	1:A:2048:ILE:HG12	1.93	0.51
1:A:1567:ASN:HA	1:A:1570:VAL:HG22	1.93	0.51
1:A:787:VAL:HA	1:A:790:VAL:HG12	1.92	0.50
1:A:2528:ARG:HG3	1:A:2853:ALA:HB3	1.94	0.50
1:A:5048:TYR:HE1	1:A:5095:PRO:HB3	1.76	0.50
1:A:1300:SER:HA	1:A:1303:VAL:HG12	1.92	0.50
1:A:3012:LEU:HD21	1:A:3062:LEU:HD21	1.92	0.50
1:A:3254:SER:OG	1:A:3404:LEU:O	2.29	0.50
1:A:2450:ALA:HB1	1:A:2459:ILE:HD13	1.94	0.50
1:A:3115:VAL:O	1:A:3119:GLN:HB2	2.10	0.50
1:A:3958:VAL:HG11	1:A:3971:ILE:HG21	1.93	0.50
1:A:4794:LEU:HD12	1:A:4798:TYR:HD2	1.75	0.50
1:A:1421:ASP:HA	1:A:1424:ARG:HE	1.77	0.50
1:A:3645:MET:HB2	1:A:3647:LEU:HD23	1.93	0.49
1:A:4547:ALA:HA	1:A:4550:VAL:HG12	1.94	0.49
1:A:1115:SER:HA	1:A:1118:LYS:HG2	1.93	0.49
1:A:4164:GLN:HG3	1:A:4169:ILE:HG22	1.93	0.49
1:A:1841:ARG:O	1:A:1845:THR:OG1	2.29	0.49
1:A:3693:PRO:HA	1:A:3696:VAL:HG22	1.93	0.49
1:A:5069:LEU:HD11	1:A:5138:ALA:HB1	1.95	0.49
1:A:942:PRO:HB3	1:A:3207:ARG:CZ	2.42	0.49
1:A:1497:LEU:HD13	1:A:1558:MET:HE1	1.94	0.49
1:A:1704:VAL:HA	1:A:1707:THR:HG22	1.93	0.49
1:A:955:CYS:O	1:A:997:LYS:NZ	2.38	0.49
1:A:4775:PHE:HE2	1:A:4812:ILE:HG22	1.78	0.49
1:A:2189:ASP:HB3	1:A:2269:PHE:HB2	1.94	0.49
1:A:3505:ARG:NE	1:A:3590:ASP:OD2	2.45	0.49
1:A:2608:VAL:HG22	1:A:2722:LEU:HG	1.93	0.49
1:A:4113:LEU:O	1:A:4168:ARG:NH1	2.45	0.49
1:A:995:LEU:HD22	1:A:1047:ILE:HG12	1.95	0.49
1:A:1322:ILE:HA	1:A:1325:VAL:HG12	1.95	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:2022:THR:HG22	1:A:2062:PRO:HG3	1.94	0.49
1:A:4835:ASN:HA	1:A:4838:VAL:HG22	1.95	0.49
1:A:3982:CYS:HB3	1:A:3987:THR:H	1.78	0.49
1:A:4271:ASP:HB3	1:A:4651:ARG:HH22	1.77	0.49
1:A:5095:PRO:HA	1:A:5128:MET:HB3	1.94	0.49
1:A:1146:HIS:NE2	1:A:2872:ILE:HG21	2.28	0.48
1:A:5033:ARG:NH2	1:A:5039:GLU:O	2.45	0.48
1:A:4165:GLU:OE2	1:A:4674:GLN:NE2	2.46	0.48
1:A:1860:ASP:O	1:A:1896:ARG:NH1	2.46	0.48
1:A:3610:LEU:HD23	1:A:3656:PRO:HD2	1.96	0.48
1:A:1031:TRP:HE3	1:A:1034:VAL:HG13	1.77	0.48
1:A:1140:GLN:HG3	1:A:1200:LEU:HD12	1.95	0.48
1:A:4816:ARG:HE	1:A:4921:THR:HG21	1.78	0.48
1:A:4953:THR:OG1	1:A:5012:ARG:NH2	2.46	0.48
1:A:3928:LEU:HD23	1:A:4069:PRO:HG3	1.95	0.48
1:A:1788:GLN:HG3	1:A:1791:GLN:HB2	1.95	0.48
1:A:2509:SER:O	1:A:2513:THR:OG1	2.31	0.48
1:A:669:PRO:HG2	1:A:723:LEU:HD23	1.96	0.48
1:A:3028:ILE:HD13	1:A:3059:VAL:HG13	1.96	0.48
1:A:3223:PRO:HA	1:A:3226:VAL:HG12	1.95	0.48
1:A:5070:ASP:OD1	1:A:5070:ASP:N	2.46	0.48
1:A:506:LEU:HD23	1:A:509:ILE:HD11	1.95	0.48
1:A:570:MET:HE1	1:A:620:LEU:HD11	1.96	0.48
1:A:1325:VAL:HG11	1:A:1381:LEU:HD12	1.94	0.48
1:A:3687:GLU:OE2	1:A:3691:LYS:NZ	2.47	0.48
1:A:3849:GLU:OE2	1:A:3867:ARG:NH2	2.44	0.48
1:A:1619:LEU:HB3	1:A:1621:LEU:HD23	1.96	0.47
1:A:3365:GLN:NE2	1:A:3369:ASP:O	2.47	0.47
1:A:4225:ARG:HH11	1:A:4256:PRO:HA	1.79	0.47
1:A:4499:THR:OG1	1:A:4500:GLN:N	2.46	0.47
1:A:2486:ASN:HB2	1:A:2529:VAL:HG11	1.96	0.47
1:A:2808:GLY:HA3	1:A:2848:TRP:HB2	1.97	0.47
1:A:3053:MET:HE1	1:A:3090:LEU:HD21	1.95	0.47
1:A:3860:ARG:HA	1:A:3863:ILE:HG12	1.96	0.47
1:A:4117:LYS:HZ3	1:A:4168:ARG:HD3	1.79	0.47
1:A:4138:ALA:HB2	1:A:4874:PRO:HG3	1.96	0.47
1:A:600:PHE:O	1:A:604:ALA:HB2	2.14	0.47
1:A:889:PHE:HB3	1:A:937:ARG:HH22	1.80	0.47
1:A:2038:LEU:HD23	1:A:2040:TYR:HE1	1.78	0.47
1:A:2369:MET:HA	1:A:2372:ARG:HB3	1.95	0.47
1:A:3144:LYS:HA	1:A:3147:VAL:HG12	1.95	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:3662:ARG:HD2	1:A:3720:LEU:HD13	1.95	0.47
1:A:4639:LEU:HA	1:A:4642:ILE:HG22	1.96	0.47
1:A:1031:TRP:CZ3	1:A:1033:ASP:HB2	2.49	0.47
1:A:3200:ASP:N	1:A:3200:ASP:OD1	2.48	0.47
1:A:2094:ARG:HB3	1:A:2098:GLU:HB2	1.96	0.47
1:A:4762:ASP:N	1:A:4762:ASP:OD1	2.45	0.47
1:A:4130:ASP:OD2	1:A:4882:TYR:OH	2.28	0.47
1:A:4896:LEU:HA	1:A:4899:ILE:HG22	1.97	0.47
1:A:4039:ILE:HA	1:A:4042:LEU:HG	1.97	0.47
1:A:1495:SER:HB3	1:A:1498:SER:H	1.80	0.46
1:A:2866:PRO:HG3	1:A:2914:LEU:HD21	1.97	0.46
1:A:1002:LEU:HD22	1:A:1037:LEU:HD12	1.97	0.46
1:A:2847:ASN:ND2	6:A:5206:AGS:S1G	2.83	0.46
1:A:3185:CYS:HA	1:A:3188:VAL:HG12	1.98	0.46
1:A:3726:SER:OG	1:A:3728:GLU:OE1	2.32	0.46
1:A:3785:HIS:HD2	1:A:3788:VAL:HB	1.79	0.46
1:A:642:ARG:HD2	1:A:4765:ARG:HH22	1.79	0.46
1:A:2728:LEU:HD12	1:A:2845:ILE:HG12	1.97	0.46
1:A:2999:SER:O	1:A:3077:GLN:NE2	2.48	0.46
1:A:1865:GLU:OE2	1:A:1869:GLN:NE2	2.48	0.46
1:A:4050:LYS:HD2	1:A:4061:ARG:HH11	1.81	0.46
1:A:626:GLY:HA2	1:A:631:LEU:HD23	1.96	0.46
1:A:720:LYS:HB3	1:A:723:LEU:HD13	1.98	0.46
1:A:1942:PHE:HD2	1:A:1946:LEU:HB2	1.79	0.46
1:A:3006:THR:HG22	1:A:3134:ASP:HA	1.96	0.46
1:A:1020:ASP:OD1	1:A:1021:VAL:N	2.44	0.46
1:A:1203:LEU:HD11	1:A:1290:MET:HG3	1.98	0.46
1:A:3014:ILE:HD11	1:A:3138:VAL:HG21	1.98	0.46
1:A:3958:VAL:HG21	1:A:3993:PHE:HB2	1.97	0.46
1:A:4396:VAL:HG11	1:A:4535:LEU:HB3	1.97	0.46
1:A:1937:SER:OG	1:A:1938:ALA:N	2.49	0.46
1:A:2406:THR:O	1:A:2443:THR:OG1	2.28	0.46
1:A:2587:MET:HE1	1:A:2658:LEU:HD11	1.98	0.46
1:A:3418:GLN:HE22	1:A:3551:THR:HG1	1.57	0.46
1:A:1986:ARG:NH1	1:A:2019:ASP:OD2	2.50	0.45
1:A:2227:THR:HA	1:A:2257:LYS:HG3	1.97	0.45
1:A:3946:PRO:HG2	1:A:4072:ASP:HB2	1.97	0.45
1:A:4126:SER:HA	1:A:4129:GLU:HG2	1.96	0.45
1:A:2182:PHE:HZ	1:A:2219:MET:HG3	1.81	0.45
1:A:3327:ASN:O	1:A:3332:LYS:NZ	2.42	0.45
1:A:2272:ASP:N	1:A:2272:ASP:OD1	2.48	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:3135:MET:HE2	1:A:3186:ALA:HB1	1.98	0.45
1:A:2303:LYS:HD2	1:A:2303:LYS:HA	1.84	0.45
1:A:4114:THR:HG23	1:A:4168:ARG:HH22	1.81	0.45
1:A:5099:LEU:HA	1:A:5128:MET:HE1	1.98	0.45
1:A:5107:MET:HG3	1:A:5111:ASP:HB3	1.99	0.45
1:A:1820:THR:OG1	1:A:1822:ASP:OD1	2.28	0.45
1:A:1981:PRO:HD2	1:A:2014:VAL:HG22	1.99	0.45
1:A:3880:VAL:HG23	1:A:3884:LEU:HD12	1.99	0.45
1:A:1456:GLN:HG3	1:A:1460:ARG:HH21	1.82	0.45
1:A:1663:LEU:HD11	1:A:1761:LEU:HD12	1.99	0.45
1:A:1286:PHE:O	1:A:1289:THR:OG1	2.35	0.45
1:A:1826:LEU:HD11	1:A:2033:PHE:HD2	1.82	0.45
1:A:2624:ASP:O	1:A:2628:HIS:HB2	2.17	0.45
1:A:646:ARG:HD3	1:A:646:ARG:HA	1.62	0.45
1:A:1082:GLN:OE1	1:A:1082:GLN:N	2.48	0.45
1:A:1828:THR:O	1:A:1831:THR:OG1	2.30	0.45
1:A:3194:GLU:HG2	1:A:3199:ARG:HH22	1.82	0.45
1:A:3519:ARG:HD2	1:A:3598:LEU:HD12	1.98	0.45
1:A:4577:MET:HG2	1:A:4693:GLU:HB3	1.99	0.45
1:A:706:LYS:NZ	1:A:707:ALA:O	2.42	0.45
1:A:849:LEU:HD23	1:A:849:LEU:HA	1.84	0.45
1:A:2307:MET:HE1	1:A:2312:PHE:HD1	1.82	0.45
1:A:2935:LEU:HB3	1:A:2940:ILE:HD11	1.98	0.45
1:A:3140:GLN:OE1	1:A:3142:TRP:NE1	2.40	0.45
1:A:3809:ASP:OD2	1:A:3851:TYR:OH	2.30	0.45
1:A:1988:THR:OG1	1:A:2457:SER:OG	2.28	0.44
1:A:2830:ILE:HG22	1:A:2831:GLU:HG3	1.99	0.44
1:A:1481:TRP:HA	1:A:1484:THR:HG22	1.98	0.44
1:A:3382:MET:HE3	1:A:5036:LEU:HD13	1.99	0.44
1:A:3904:LEU:HD23	1:A:3907:LYS:HZ3	1.82	0.44
1:A:4046:LEU:HD11	1:A:4084:LEU:HD12	1.99	0.44
1:A:4167:VAL:HG23	1:A:4168:ARG:HG2	1.98	0.44
1:A:5002:GLN:HG2	1:A:5003:THR:H	1.82	0.44
1:A:4201:LEU:HB3	1:A:4205:LYS:HB2	1.98	0.44
1:A:4706:GLU:HG2	1:A:4707:THR:HG23	1.98	0.44
1:A:896:ILE:HD11	1:A:899:ALA:HA	2.00	0.44
1:A:1207:HIS:NE2	1:A:2887:ASP:OD2	2.49	0.44
1:A:4753:ARG:HA	1:A:4756:ILE:HG22	1.99	0.44
1:A:2938:GLN:O	1:A:2942:HIS:ND1	2.46	0.44
1:A:1619:LEU:HD21	1:A:1755:LEU:HD13	2.00	0.44
1:A:1733:GLU:HG2	1:A:1738:GLY:HA3	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1807:LEU:HD21	1:A:1925:LEU:HD21	1.99	0.44
1:A:5080:ILE:HA	1:A:5084:LEU:HD13	2.00	0.44
1:A:5107:MET:HE3	1:A:5115:LEU:HD23	2.00	0.44
1:A:3481:VAL:O	1:A:3485:VAL:HG13	2.18	0.44
1:A:3508:ILE:HG21	1:A:3595:LEU:HB3	1.99	0.44
1:A:1018:VAL:HG23	1:A:1019:PHE:H	1.82	0.44
1:A:1640:ARG:NH1	1:A:1643:GLU:OE1	2.51	0.43
1:A:2326:PHE:HA	1:A:2329:LEU:HD13	2.00	0.43
1:A:2815:LYS:HB2	1:A:2815:LYS:HE2	1.87	0.43
1:A:802:GLN:HB3	1:A:804:ARG:H	1.82	0.43
1:A:1652:VAL:HA	1:A:1655:GLN:HG2	1.99	0.43
1:A:3382:MET:HE3	1:A:5036:LEU:HD12	1.99	0.43
1:A:3877:ALA:HA	1:A:3880:VAL:HG12	2.00	0.43
1:A:4421:MET:SD	1:A:4424:ALA:HB3	2.58	0.43
1:A:4987:MET:HE2	1:A:4987:MET:HB2	1.94	0.43
1:A:2119:GLU:OE2	2:A:5201:ATP:O2'	2.36	0.43
1:A:2725:PRO:HB3	1:A:2842:PHE:HB3	2.01	0.43
1:A:2832:ASP:OD1	1:A:2832:ASP:N	2.51	0.43
1:A:3787:GLN:O	1:A:3790:SER:OG	2.31	0.43
1:A:5113:ASP:HB3	1:A:5114:ILE:H	1.58	0.43
1:A:1384:LEU:HD12	1:A:1384:LEU:HA	1.88	0.43
1:A:4523:PRO:HG3	1:A:4696:GLN:HE21	1.83	0.43
1:A:1987:LEU:HD23	1:A:1992:LEU:HD11	2.00	0.43
1:A:3611:TRP:HD1	1:A:3656:PRO:HB3	1.83	0.43
1:A:4695:LEU:HA	1:A:4698:ILE:HG22	2.01	0.43
1:A:789:ASN:O	1:A:793:MET:HG2	2.19	0.43
1:A:2137:GLN:HG2	1:A:2139:GLN:HE22	1.83	0.43
1:A:2177:ARG:HH21	2:A:5201:ATP:H4'	1.82	0.43
1:A:4883:GLN:HB3	1:A:4891:SER:HG	1.84	0.43
1:A:1810:TYR:OH	1:A:1853:VAL:O	2.37	0.42
1:A:2035:LEU:HD11	1:A:2057:TYR:HE2	1.84	0.42
1:A:2632:ASP:OD1	1:A:2632:ASP:N	2.48	0.42
1:A:2096:PRO:HG3	1:A:2173:TRP:CE2	2.54	0.42
1:A:2717:VAL:O	1:A:2721:GLU:HG2	2.19	0.42
1:A:1956:ALA:HB3	1:A:2534:PRO:HB2	2.01	0.42
1:A:2081:LEU:HB3	1:A:2085:LEU:HD23	2.00	0.42
1:A:2447:PHE:HB3	1:A:2450:ALA:HB2	2.01	0.42
1:A:4291:VAL:HG21	1:A:4324:GLU:HB3	2.01	0.42
1:A:4680:CYS:SG	1:A:4681:SER:N	2.92	0.42
1:A:1552:LEU:O	1:A:1556:MET:HB2	2.20	0.42
1:A:3790:SER:O	1:A:3793:SER:HB2	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:4677:ILE:HG23	1:A:4678:ILE:HD12	2.01	0.42
1:A:646:ARG:HH12	1:A:675:MET:HB3	1.84	0.42
1:A:2041:LEU:HB2	1:A:2049:TRP:HB3	2.01	0.42
1:A:3855:SER:O	1:A:3860:ARG:NH1	2.52	0.42
1:A:668:LEU:O	1:A:672:HIS:ND1	2.53	0.42
1:A:1822:ASP:OD1	1:A:1822:ASP:N	2.53	0.42
1:A:2130:GLN:HE22	1:A:2151:VAL:HB	1.85	0.42
1:A:2473:LEU:HD23	1:A:2473:LEU:HA	1.89	0.42
1:A:3516:ASP:OD1	1:A:3516:ASP:N	2.52	0.42
1:A:5045:ASP:OD2	1:A:5047:GLN:NE2	2.52	0.42
1:A:685:LYS:HE2	1:A:685:LYS:HB3	1.82	0.42
1:A:3000:ARG:NH1	1:A:3126:ASN:O	2.53	0.42
1:A:2984:TYR:OH	1:A:3021:GLU:O	2.36	0.42
1:A:3878:LEU:HD22	1:A:4026:MET:HE1	2.02	0.42
1:A:1528:LEU:HD23	1:A:1548:LEU:HD21	2.01	0.41
1:A:3592:ASP:OD1	1:A:3592:ASP:N	2.53	0.41
1:A:1161:GLN:OE1	1:A:1165:ASN:ND2	2.50	0.41
1:A:1190:SER:HB2	1:A:1192:ASP:OD1	2.21	0.41
1:A:600:PHE:O	1:A:604:ALA:CB	2.69	0.41
1:A:792:LYS:HA	1:A:795:MET:HG2	2.02	0.41
1:A:3274:PHE:HD1	1:A:3373:VAL:HB	1.85	0.41
1:A:690:ASP:HB3	1:A:696:GLU:HA	2.01	0.41
1:A:2208:PHE:HD1	1:A:2311:LEU:HD21	1.85	0.41
1:A:3224:ASP:OD1	1:A:3224:ASP:N	2.51	0.41
1:A:3612:MET:HE1	1:A:3615:TYR:HD2	1.85	0.41
1:A:4098:ASP:N	1:A:4098:ASP:OD1	2.53	0.41
1:A:4129:GLU:HA	1:A:4132:VAL:HG12	2.01	0.41
1:A:1351:PHE:HD2	1:A:1353:ASN:H	1.67	0.41
1:A:3068:TYR:HB2	1:A:3124:LEU:HD11	2.03	0.41
1:A:521:LEU:HA	1:A:524:PHE:HD2	1.86	0.41
1:A:1120:ARG:HD2	1:A:1167:LYS:NZ	2.36	0.41
1:A:1319:ALA:HA	1:A:1322:ILE:HG22	2.01	0.41
1:A:1882:HIS:NE2	1:A:1884:GLU:OE1	2.50	0.41
1:A:3200:ASP:H	1:A:3205:LEU:HD13	1.84	0.41
1:A:937:ARG:NH1	1:A:3237:PHE:CE1	2.89	0.41
1:A:1649:TRP:HA	1:A:1652:VAL:HG12	2.01	0.41
1:A:3736:ALA:HB2	1:A:3808:LEU:HD12	2.02	0.41
1:A:4562:PHE:HD2	1:A:4563:LEU:HD22	1.85	0.41
1:A:4580:ARG:HB3	1:A:4584:GLU:HB2	2.02	0.41
1:A:1283:ASP:O	1:A:1287:MET:HB2	2.21	0.41
1:A:2556:GLN:HA	1:A:2559:VAL:HG12	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:3558:ASP:OD2	1:A:5042:ARG:NH1	2.53	0.41
1:A:3726:SER:OG	1:A:3727:ARG:N	2.54	0.41
1:A:3821:LYS:HZ1	1:A:3872:ARG:HH22	1.68	0.41
1:A:3960:LEU:HG	1:A:3989:LEU:HD12	2.03	0.41
1:A:4534:HIS:HA	1:A:4537:MET:HG3	2.03	0.41
1:A:4722:ARG:O	1:A:4725:LYS:NZ	2.46	0.41
1:A:942:PRO:HA	1:A:943:PRO:HD3	1.91	0.41
1:A:2177:ARG:NE	2:A:5201:ATP:O3'	2.53	0.41
1:A:3133:LEU:HD11	1:A:3138:VAL:HB	2.03	0.41
1:A:4338:PRO:HA	1:A:4342:GLN:HB2	2.02	0.41
1:A:4698:ILE:O	1:A:4702:LYS:HG2	2.21	0.41
1:A:4883:GLN:HB3	1:A:4891:SER:OG	2.21	0.41
1:A:1005:VAL:O	1:A:1009:SER:HB3	2.21	0.40
1:A:1635:LEU:HD13	1:A:1638:LEU:HD23	2.03	0.40
1:A:1697:LYS:HD2	1:A:1697:LYS:HA	1.89	0.40
1:A:2742:ILE:O	1:A:2746:ALA:CB	2.69	0.40
1:A:3015:LEU:HD23	1:A:3015:LEU:HA	1.93	0.40
1:A:1494:LEU:O	1:A:1495:SER:HB2	2.22	0.40
1:A:1010:TRP:HE1	1:A:1018:VAL:HG11	1.86	0.40
1:A:1120:ARG:HD2	1:A:1167:LYS:HZ1	1.86	0.40
1:A:1401:ILE:HD12	1:A:1404:LEU:HD12	2.02	0.40
1:A:3282:ARG:HD2	1:A:3282:ARG:HA	1.83	0.40
1:A:4009:ALA:HA	1:A:4012:ARG:HG2	2.03	0.40
1:A:3549:TRP:CD1	1:A:3549:TRP:H	2.40	0.40
1:A:4259:VAL:HG23	1:A:4656:PRO:HB2	2.04	0.40
1:A:4281:LYS:HA	1:A:4284:ARG:HG2	2.04	0.40
1:A:4580:ARG:HB2	1:A:4585:THR:HG23	2.04	0.40
1:A:4711:LEU:HD21	1:A:4899:ILE:HG23	2.03	0.40
1:A:5024:PHE:HA	1:A:5074:LEU:HD11	2.04	0.40
1:A:5100:LYS:HE2	1:A:5119:GLU:HA	2.04	0.40
1:A:938:LEU:HD11	1:A:968:PHE:CE1	2.56	0.40
1:A:1588:ILE:HD12	1:A:1588:ILE:HA	1.88	0.40
1:A:3922:VAL:HA	1:A:3925:MET:HG2	2.02	0.40
1:A:3975:LEU:HD23	1:A:3993:PHE:HD1	1.86	0.40
1:A:4925:ARG:HH22	1:A:4929:ASN:HB3	1.86	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	4394/4836 (91%)	4028 (92%)	358 (8%)	8 (0%)	43 75

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	3122	VAL
1	A	4600	GLN
1	A	1374	SER
1	A	3496	ASN
1	A	4166	PRO
1	A	5128	MET
1	A	4049	GLN
1	A	1373	ILE

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	3970/4342 (91%)	3968 (100%)	2 (0%)	88 89

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4587	HIS
1	A	4842	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (34) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	500	GLN
1	A	1255	ASN
1	A	1282	ASN
1	A	1378	GLN
1	A	1395	HIS
1	A	1467	HIS
1	A	1595	ASN
1	A	1799	HIS
1	A	2045	ASN
1	A	2270	ASN
1	A	2557	GLN
1	A	2655	HIS
1	A	2677	ASN
1	A	2702	ASN
1	A	2764	HIS
1	A	2777	GLN
1	A	2790	GLN
1	A	2953	ASN
1	A	3095	HIS
1	A	3267	HIS
1	A	3419	ASN
1	A	3734	GLN
1	A	3838	ASN
1	A	4235	GLN
1	A	4263	GLN
1	A	4337	HIS
1	A	4342	GLN
1	A	4404	HIS
1	A	4560	GLN
1	A	4564	GLN
1	A	4696	GLN
1	A	4716	GLN
1	A	4885	GLN
1	A	5121	GLN

5.3.3 RNA ⓘ

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 6 ligands modelled in this entry, 3 are monoatomic - leaving 3 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
5	ADP	A	5205	-	28,29,29	1.42	4 (14%)	43,45,45	1.90	9 (20%)
6	AGS	A	5206	-	32,33,33	0.51	1 (3%)	45,52,52	0.70	1 (2%)
2	ATP	A	5201	3	32,33,33	0.37	0	48,52,52	0.33	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	ADP	A	5205	-	-	5/16/32/32	0/3/3/3
6	AGS	A	5206	-	-	8/21/38/38	0/3/3/3
2	ATP	A	5201	3	-	7/22/38/38	0/3/3/3

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	5205	ADP	C5-C4	4.64	1.47	1.39
5	A	5205	ADP	C5-C6	2.74	1.48	1.41
5	A	5205	ADP	C8-N7	2.42	1.36	1.31

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	5205	ADP	C5-N7	-2.31	1.34	1.39
6	A	5206	AGS	PG-S1G	2.10	1.95	1.90

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	5205	ADP	C5-C4-N3	-6.15	118.25	126.72
5	A	5205	ADP	N3-C4-N9	4.79	135.31	127.17
5	A	5205	ADP	C2-N3-C4	3.92	121.41	111.83
6	A	5206	AGS	PB-O3B-PG	-3.68	119.72	133.17
5	A	5205	ADP	C4-C5-N7	-3.56	106.51	110.58
5	A	5205	ADP	N3-C2-N1	-3.40	123.43	128.58
5	A	5205	ADP	O4'-C1'-N9	2.84	113.53	108.09
5	A	5205	ADP	C5-N7-C8	2.62	107.56	103.45
5	A	5205	ADP	C4-N9-C8	2.46	108.33	105.74
5	A	5205	ADP	C6-C5-N7	2.19	136.31	132.09

There are no chirality outliers.

All (20) torsion outliers are listed below:

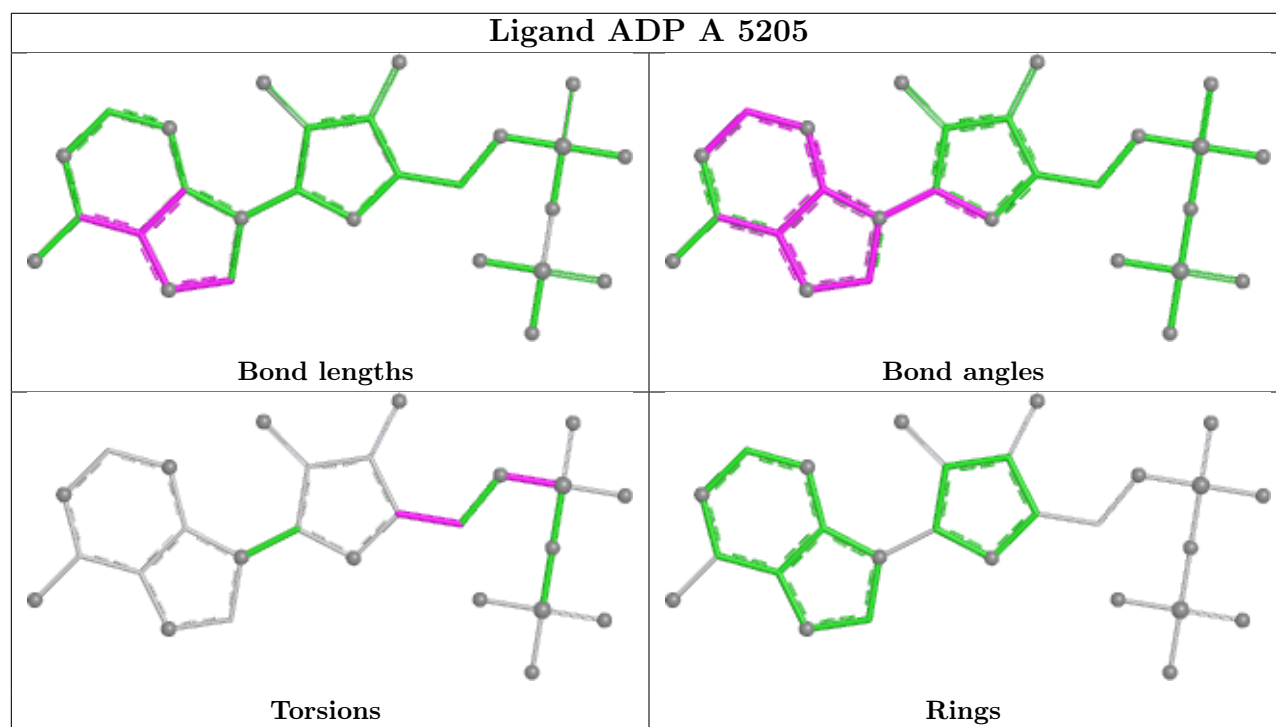
Mol	Chain	Res	Type	Atoms
2	A	5201	ATP	C5'-O5'-PA-O1A
2	A	5201	ATP	C5'-O5'-PA-O2A
2	A	5201	ATP	C5'-O5'-PA-O3A
2	A	5201	ATP	O4'-C4'-C5'-O5'
5	A	5205	ADP	C5'-O5'-PA-O2A
5	A	5205	ADP	O4'-C4'-C5'-O5'
6	A	5206	AGS	C5'-O5'-PA-O2A
6	A	5206	AGS	C5'-O5'-PA-O3A
6	A	5206	AGS	C4'-C5'-O5'-PA
5	A	5205	ADP	C3'-C4'-C5'-O5'
2	A	5201	ATP	C3'-C4'-C5'-O5'
6	A	5206	AGS	C2'-C1'-N9-C8
5	A	5205	ADP	C5'-O5'-PA-O1A
5	A	5205	ADP	C5'-O5'-PA-O3A
2	A	5201	ATP	PA-O3A-PB-O2B
6	A	5206	AGS	O4'-C1'-N9-C4
6	A	5206	AGS	O4'-C1'-N9-C8
6	A	5206	AGS	C2'-C1'-N9-C4
6	A	5206	AGS	PG-O3B-PB-O2B
2	A	5201	ATP	PA-O3A-PB-O1B

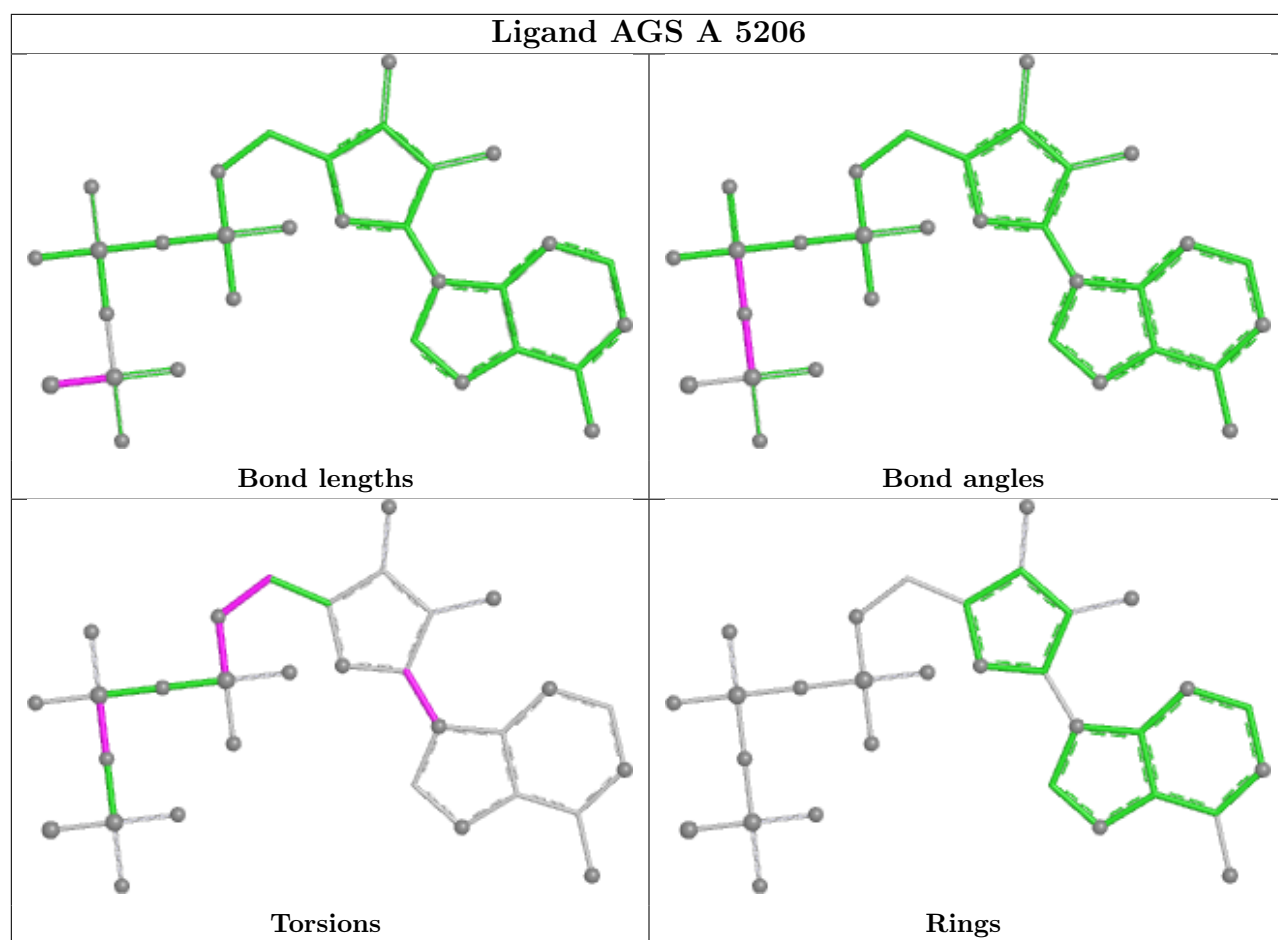
There are no ring outliers.

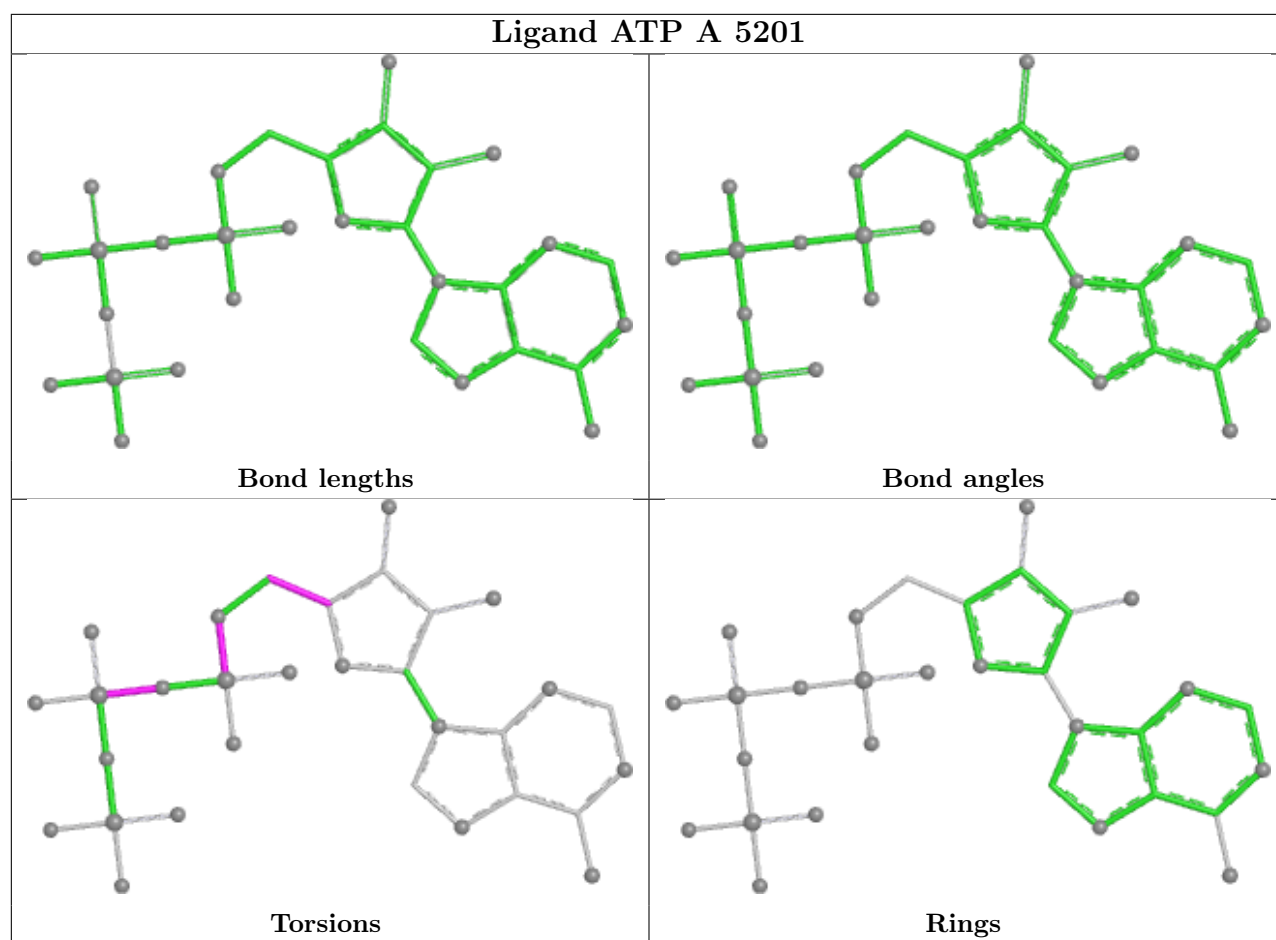
2 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	5206	AGS	2	0
2	A	5201	ATP	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 than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.







5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

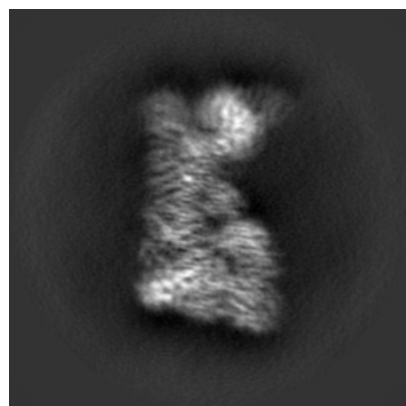
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-12932. These allow visual inspection of the internal detail of the map and identification of artifacts.

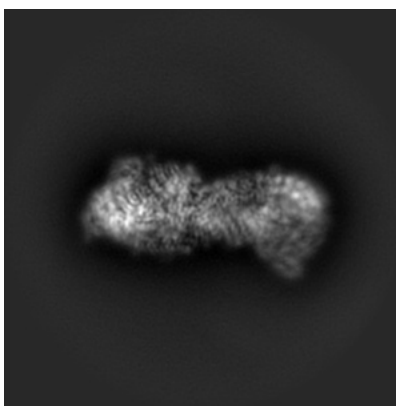
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

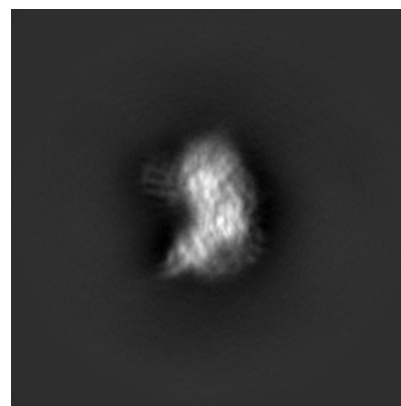
6.1.1 Primary map



X

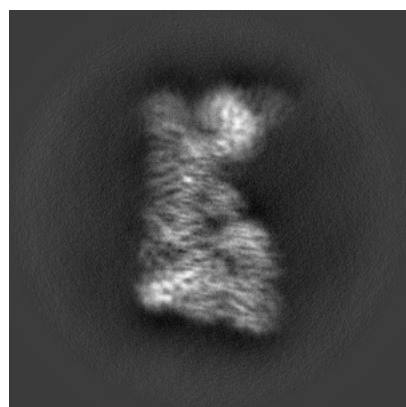


Y

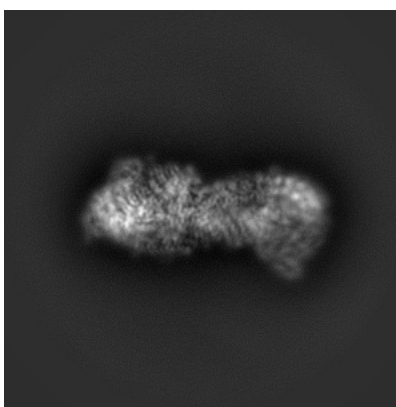


Z

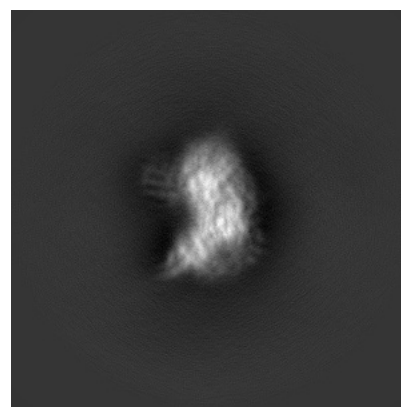
6.1.2 Raw map



X



Y

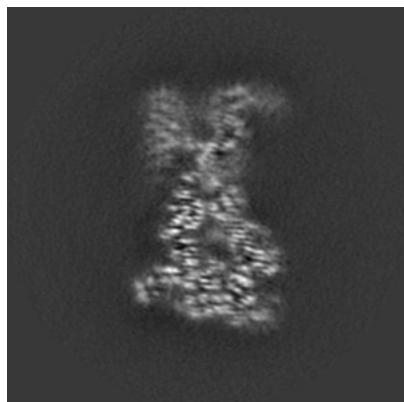


Z

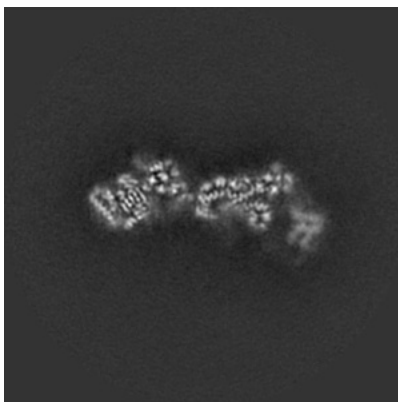
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

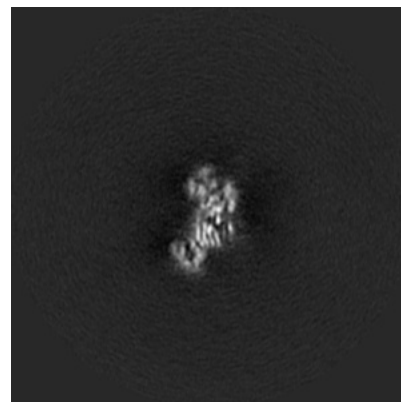
6.2.1 Primary map



X Index: 176

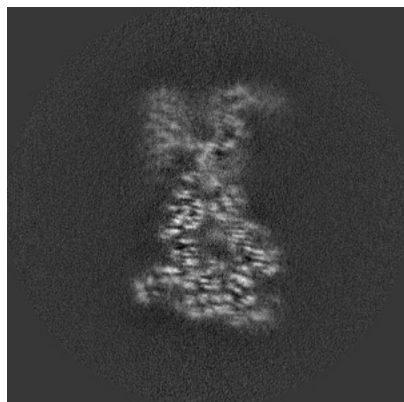


Y Index: 176

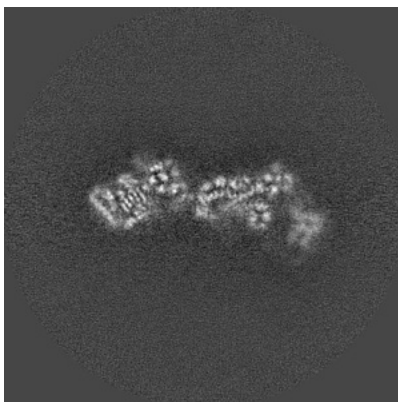


Z Index: 176

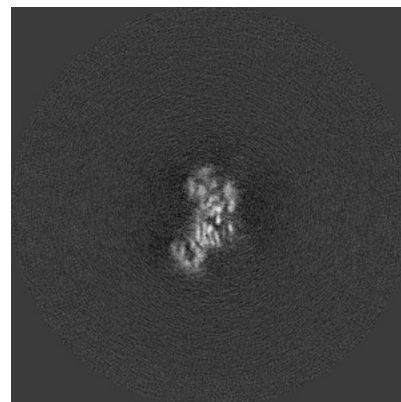
6.2.2 Raw map



X Index: 176



Y Index: 176

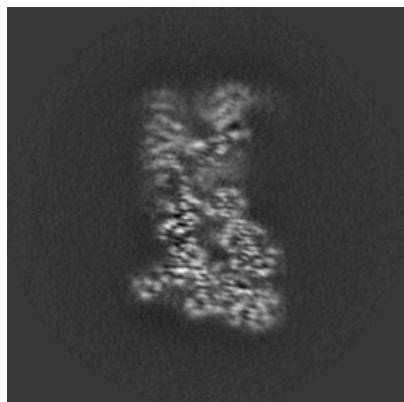


Z Index: 176

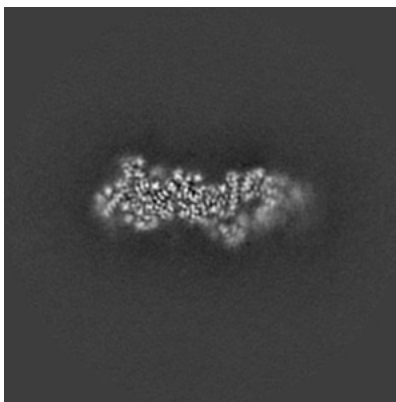
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

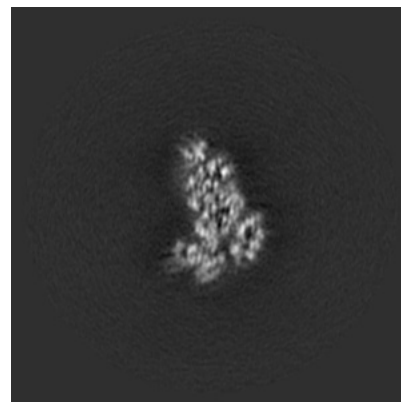
6.3.1 Primary map



X Index: 172

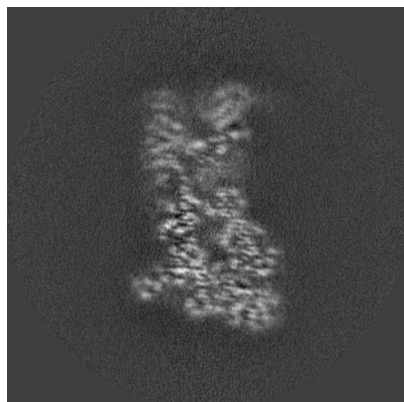


Y Index: 159

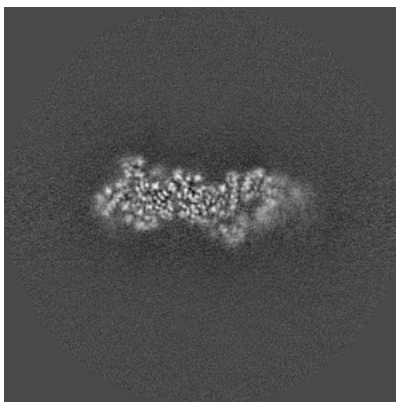


Z Index: 109

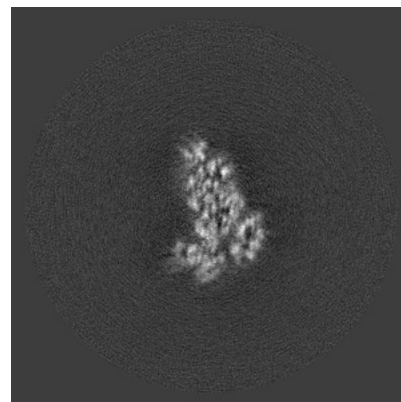
6.3.2 Raw map



X Index: 172



Y Index: 159



Z Index: 109

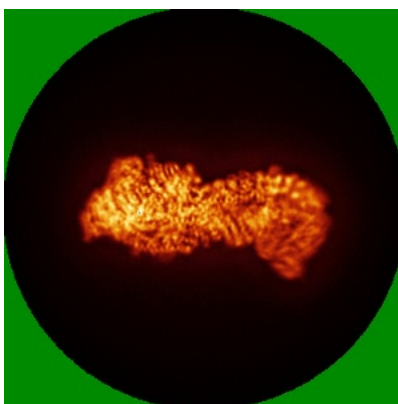
The images above show the largest variance slices of the map in three orthogonal directions.

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

6.4.1 Primary map



X

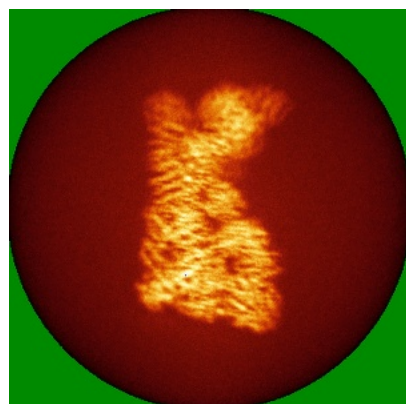


Y

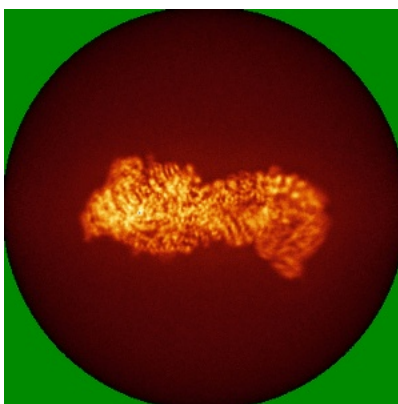


Z

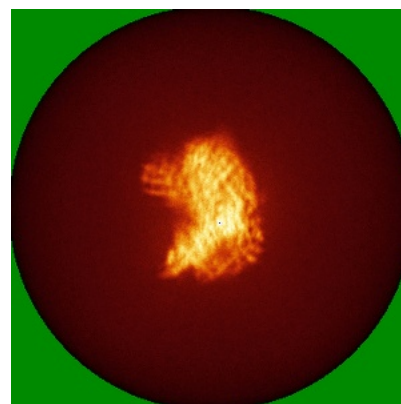
6.4.2 Raw map



X



Y

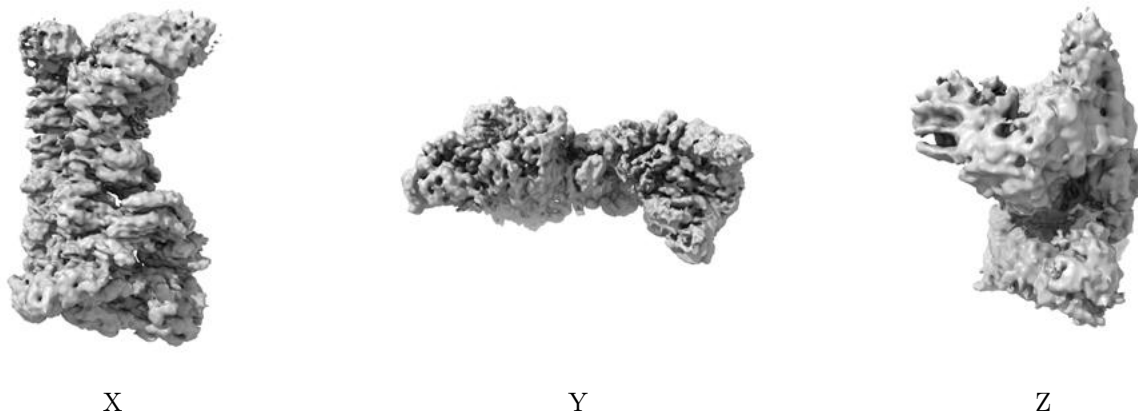


Z

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

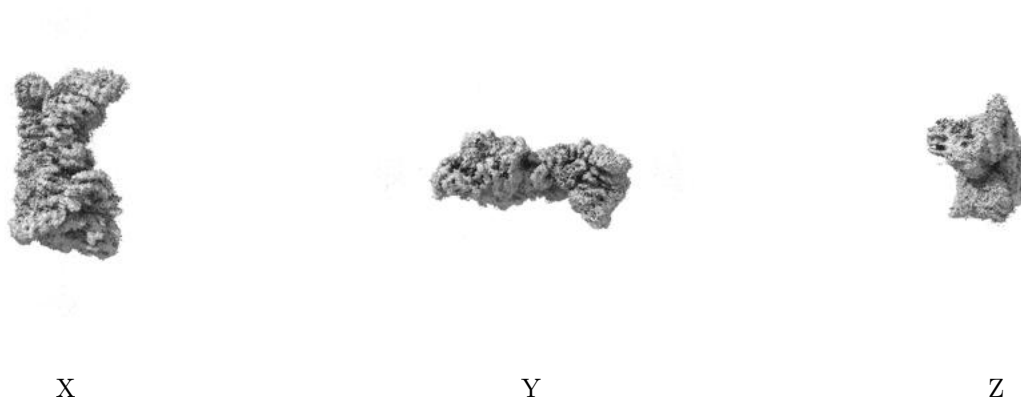
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.005. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.5.2 Raw map



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

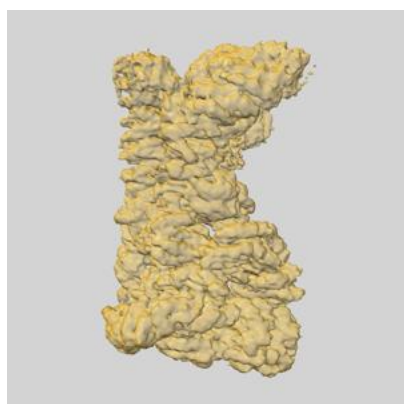
6.6 Mask visualisation [i](#)

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

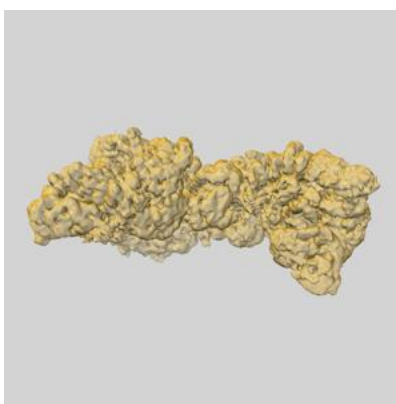
A mask typically either:

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

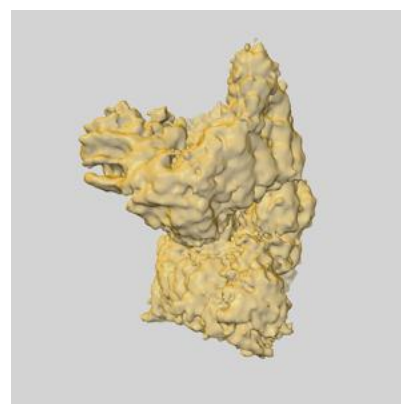
6.6.1 emd_12932_msk_2.map [i](#)



X

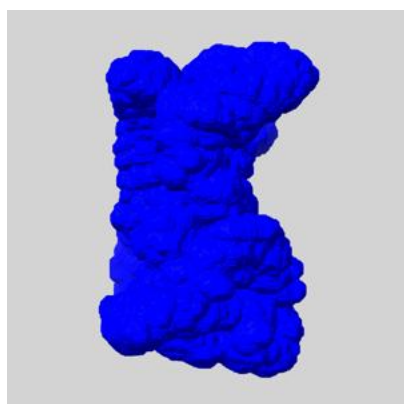


Y

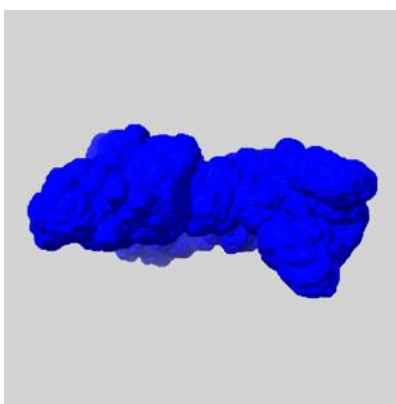


Z

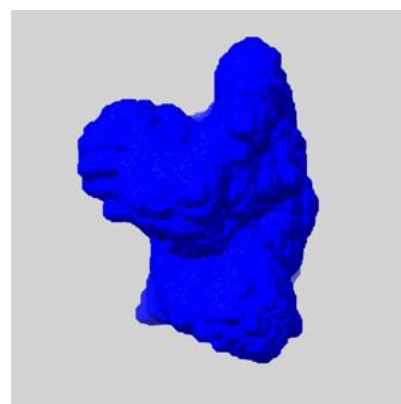
6.6.2 emd_12932_msk_1.map [i](#)



X



Y

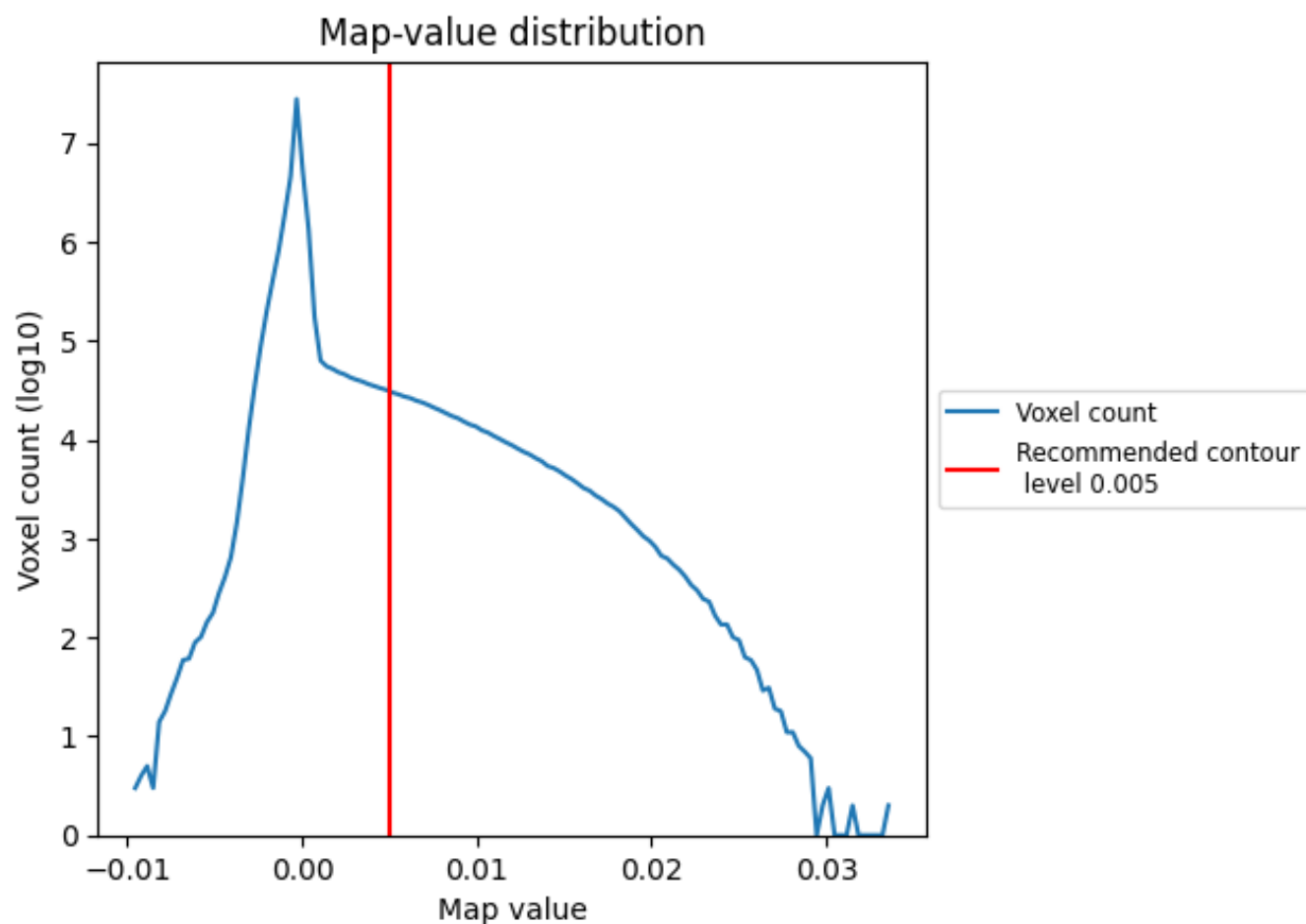


Z

7 Map analysis [i](#)

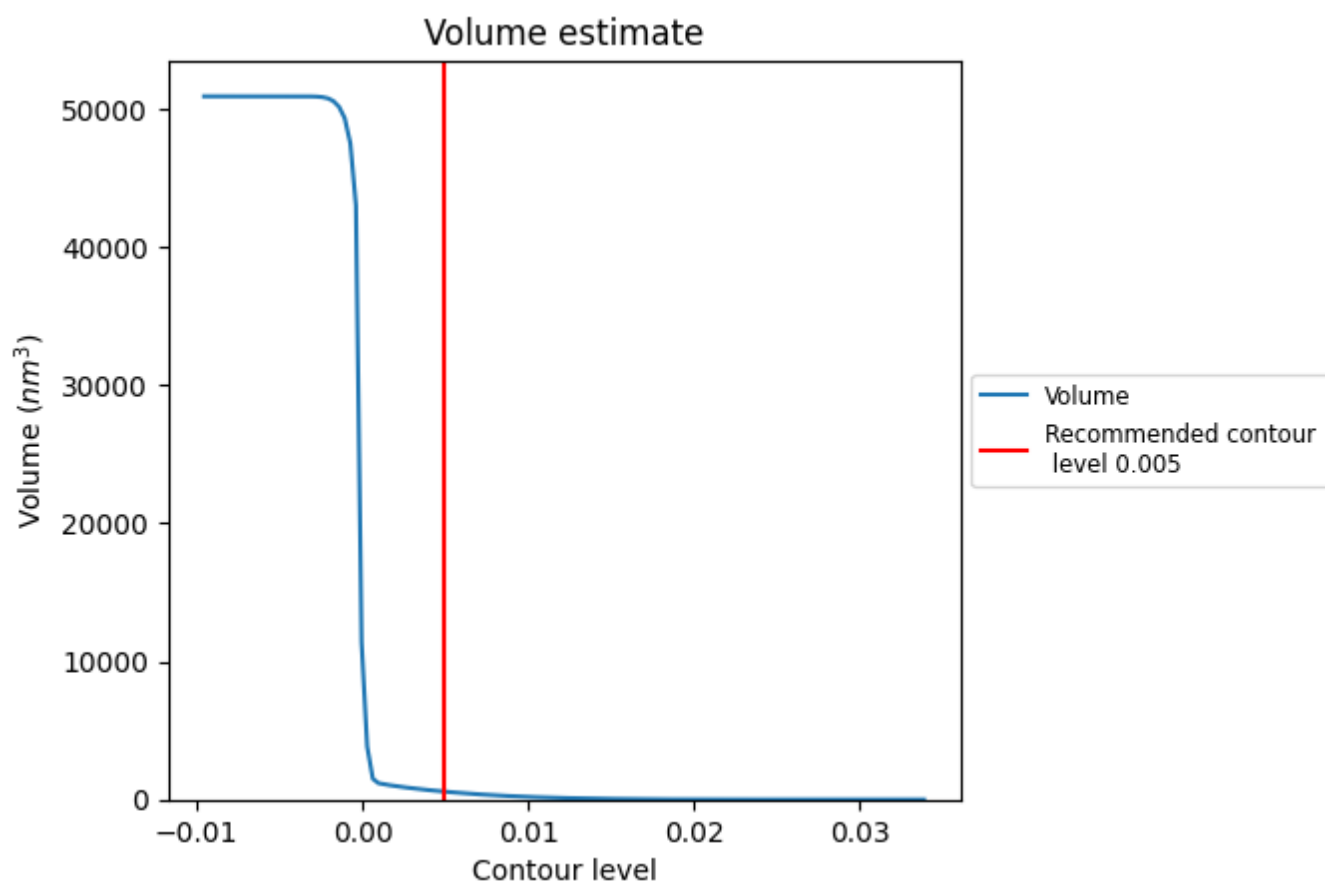
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



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

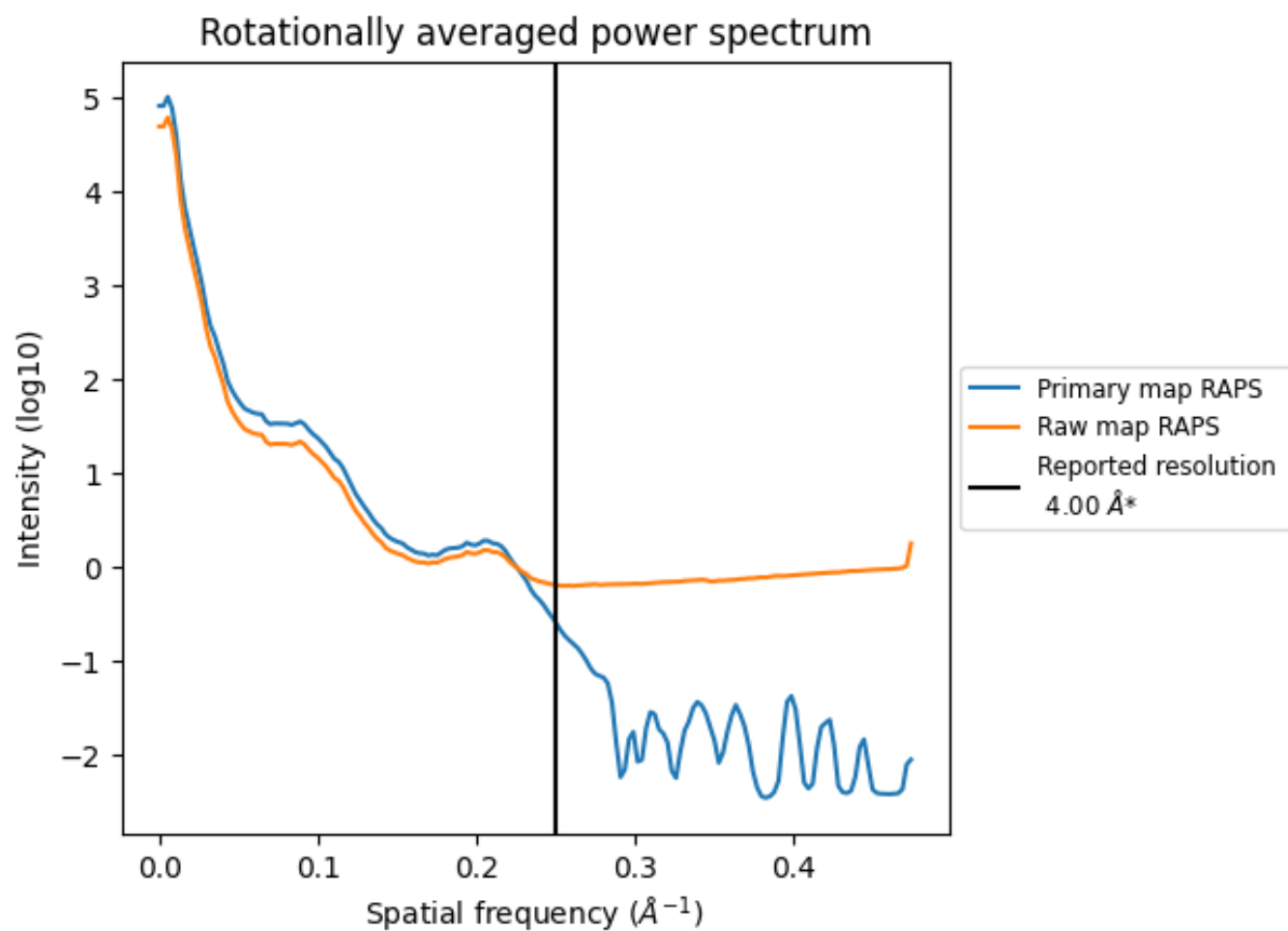
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 572 nm³; this corresponds to an approximate mass of 517 kDa.

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

7.3 Rotationally averaged power spectrum ⓘ

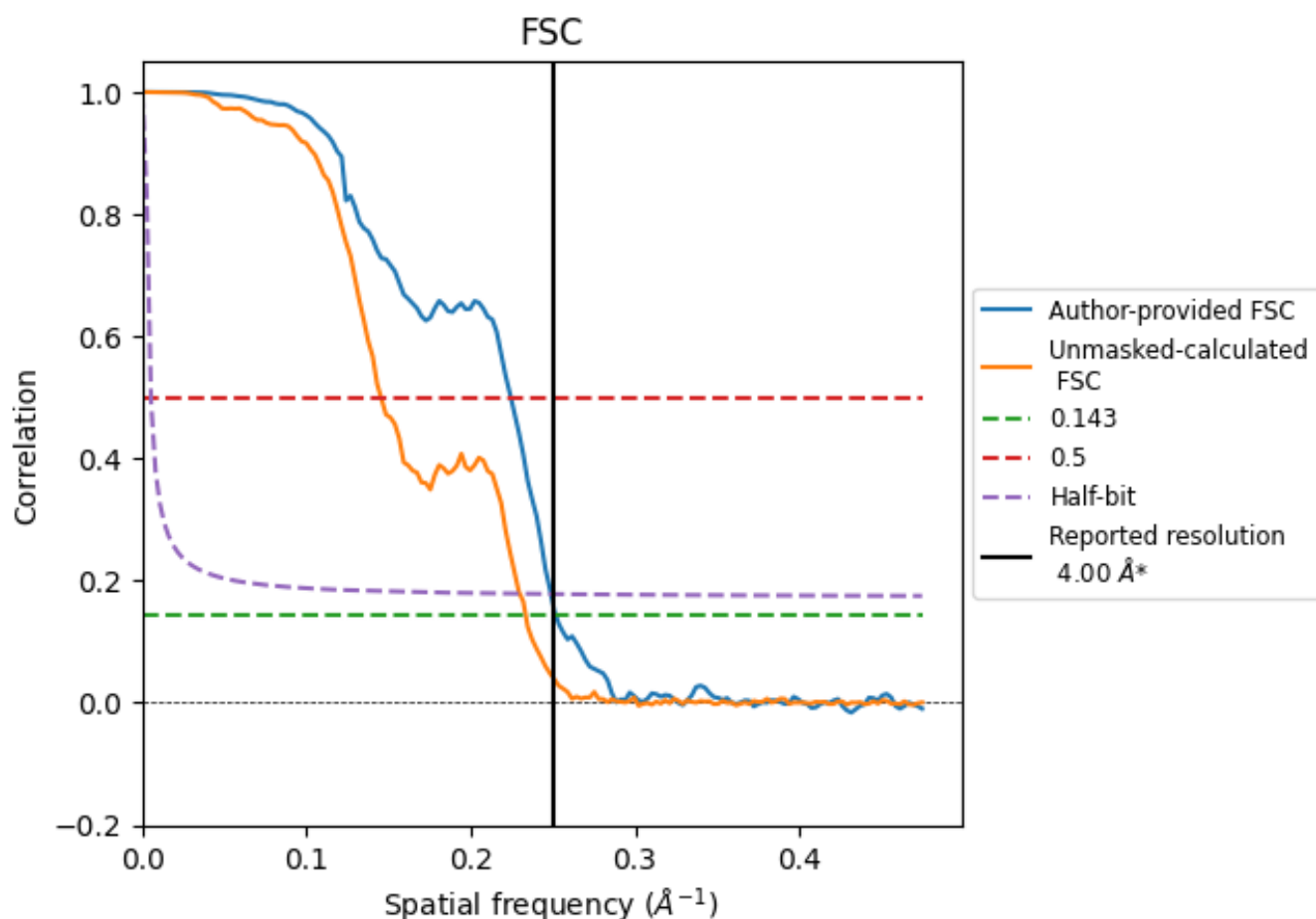


*Reported resolution corresponds to spatial frequency of 0.250 $\text{\AA}^{-1}$

8 Fourier-Shell correlation [i](#)

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

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of 0.250 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

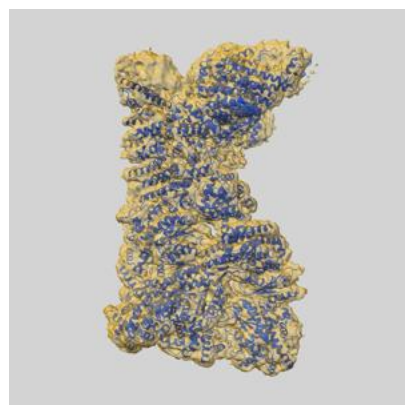
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	4.00	-	-
Author-provided FSC curve	3.97	4.46	4.02
Unmasked-calculated*	4.28	6.88	4.35

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps.

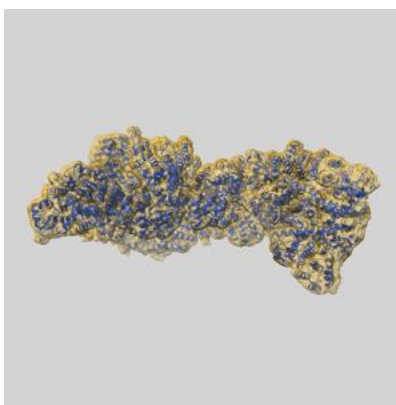
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-12932 and PDB model 7OIM. Per-residue inclusion information can be found in [section 3](#) on [page 6](#).

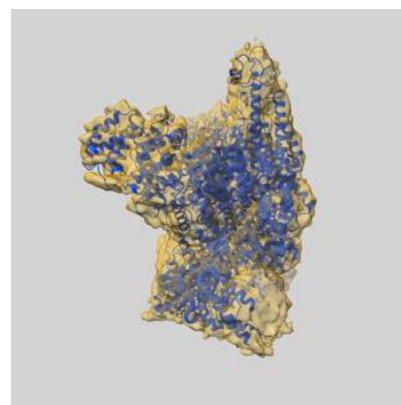
9.1 Map-model overlay [i](#)



X



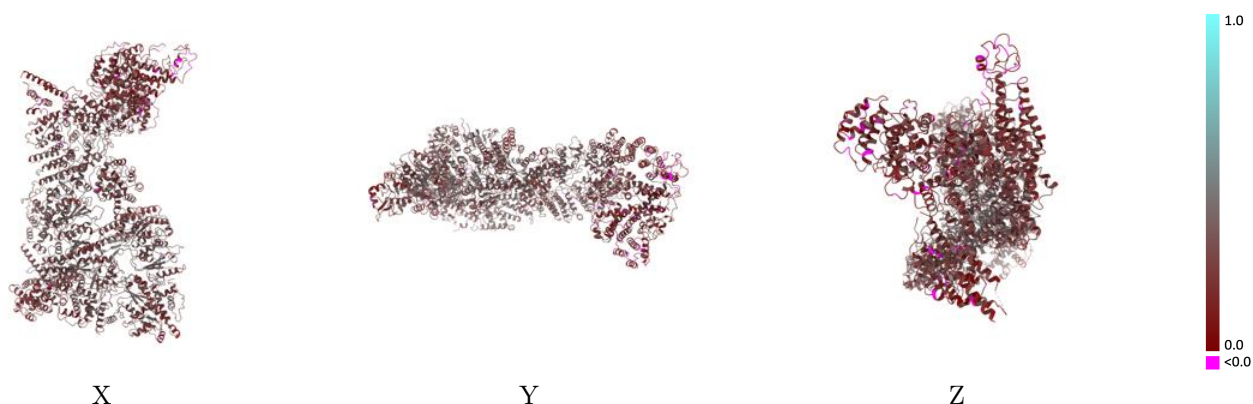
Y



Z

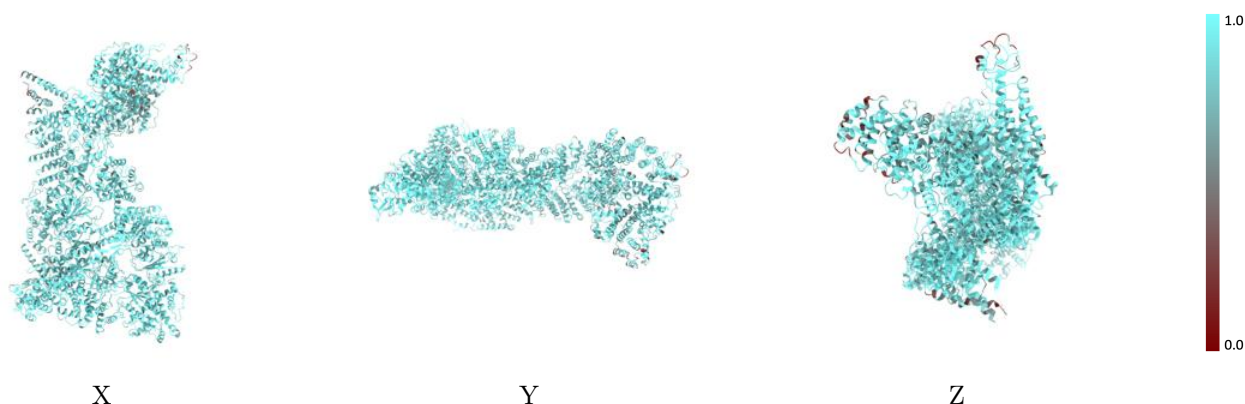
The images above show the 3D surface view of the map at the recommended contour level 0.005 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



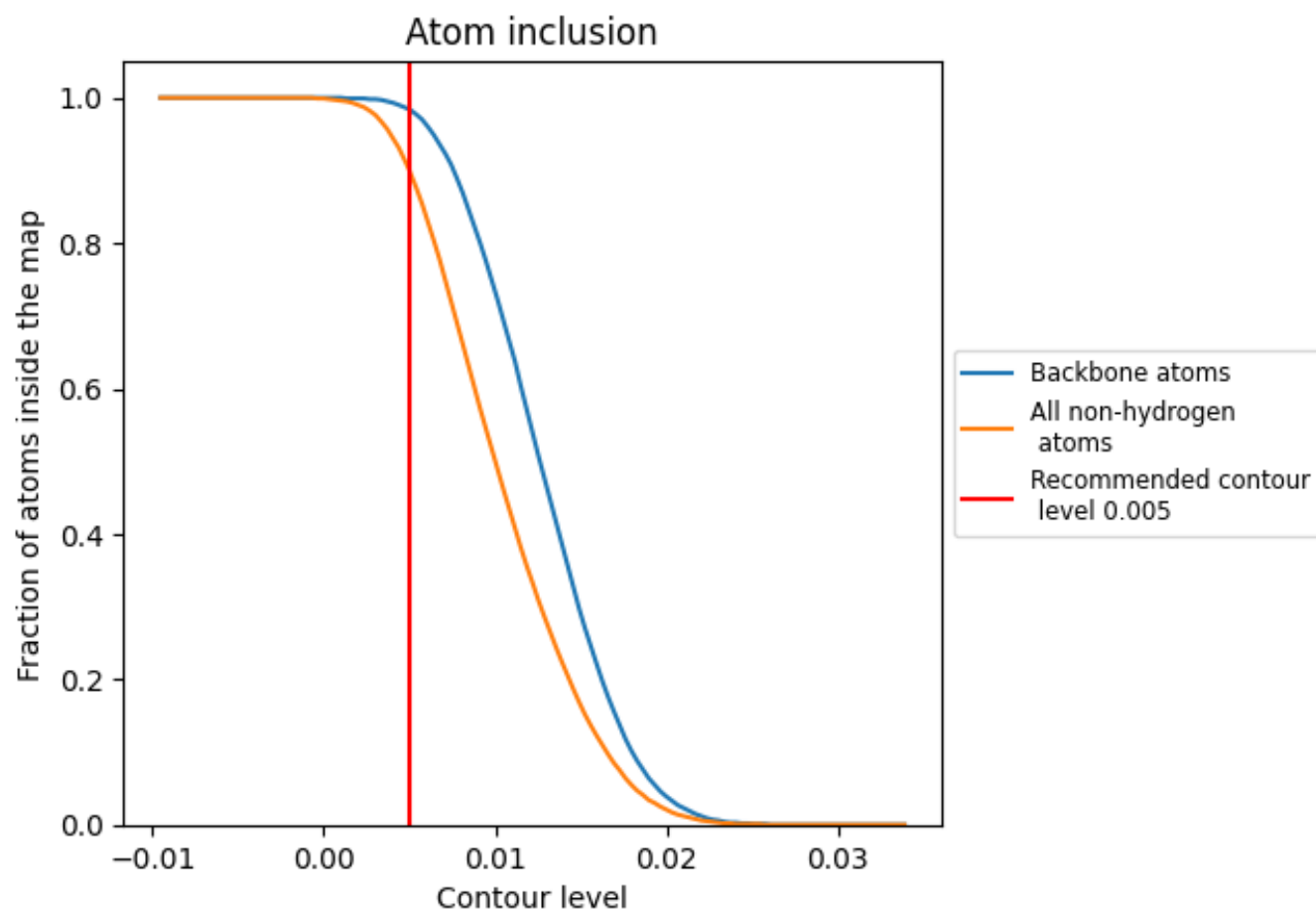
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions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

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



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (0.005).

9.4 Atom inclusion [i](#)



At the recommended contour level, 98% of all backbone atoms, 90% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary ⓘ

The table lists the average atom inclusion at the recommended contour level (0.005) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.9000	0.2870
A	0.9000	0.2870

