



Full wwPDB EM Validation Report ⓘ

Mar 29, 2026 – 05:33 PM UTC

PDB ID : 7XJJ / pdb_00007xjj
EMDB ID : EMD-33229
Title : Cryo-EM structure of the galanin-bound GALR1-miniGo complex
Authors : Jiang, W.; Zheng, S.
Deposited on : 2022-04-18
Resolution : 3.30 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

EMDB validation analysis : 0.0.1.dev132
MolProbity : 4-5-2 with Phenix2.0
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

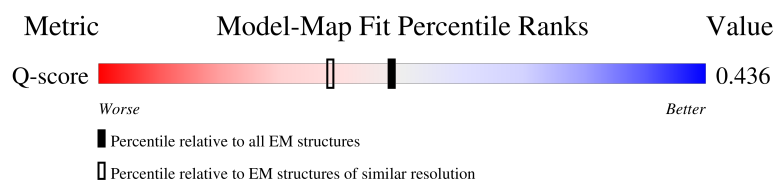
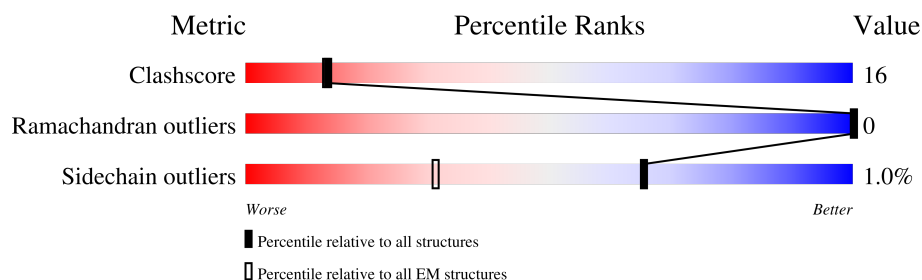
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.30 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	15087 (2.80 - 3.80)

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	228	14% 65% 26% 6%
2	B	340	19% 70% 29% .
3	C	30	7% 40% 17% 43%
4	E	347	5% 54% 21% 22%

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Mol	Chain	Length	Quality of chain
5	G	71	
6	S	298	

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 8710 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 G protein subunit alpha o1, Guanine nucleotide-binding protein G(o) subunit alpha.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	215	Total	C	N	O	S	0	0
			1691	1073	283	324	11		

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	42	ASP	GLY	conflict	UNP A0A1W2PRJ7
A	43	ASN	GLU	conflict	UNP A0A1W2PRJ7
A	174	GLY	-	linker	UNP A0A1W2PRJ7
A	175	GLY	-	linker	UNP A0A1W2PRJ7
A	176	SER	-	linker	UNP A0A1W2PRJ7
A	177	GLY	-	linker	UNP A0A1W2PRJ7
A	178	GLY	-	linker	UNP A0A1W2PRJ7
A	179	SER	-	linker	UNP A0A1W2PRJ7
A	180	GLY	-	linker	UNP A0A1W2PRJ7
A	181	GLY	-	linker	UNP A0A1W2PRJ7
A	227	ASP	ALA	conflict	UNP A0A1W2PP38
A	230	ASP	GLY	conflict	UNP A0A1W2PP38
A	332	ALA	ILE	conflict	UNP P09471
A	335	ILE	VAL	conflict	UNP P09471

- Molecule 2 is a protein called Guanine nucleotide-binding protein G(I)/G(S)/G(T) subunit beta-1.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	338	Total	C	N	O	S	0	0
			2597	1603	467	506	21		

- Molecule 3 is a protein called Galanin.

Mol	Chain	Residues	Atoms				AltConf	Trace
3	C	17	Total	C	N	O	0	0
			121	79	21	21		

- Molecule 4 is a protein called Galanin receptor type 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	E	269	Total	C	N	O	S	0	0
			2095	1407	332	344	12		

There are 25 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	-9	ASP	-	expression tag	UNP P47211
E	-8	TYR	-	expression tag	UNP P47211
E	-7	LYS	-	expression tag	UNP P47211
E	-6	ASP	-	expression tag	UNP P47211
E	-5	ASP	-	expression tag	UNP P47211
E	-4	ASP	-	expression tag	UNP P47211
E	-3	ASP	-	expression tag	UNP P47211
E	-2	LYS	-	expression tag	UNP P47211
E	-1	GLY	-	expression tag	UNP P47211
E	0	SER	-	expression tag	UNP P47211
E	323	GLY	-	expression tag	UNP P47211
E	324	GLY	-	expression tag	UNP P47211
E	325	GLY	-	expression tag	UNP P47211
E	326	GLY	-	expression tag	UNP P47211
E	327	GLY	-	expression tag	UNP P47211
E	328	GLY	-	expression tag	UNP P47211
E	329	ALA	-	expression tag	UNP P47211
E	330	GLY	-	expression tag	UNP P47211
E	331	ALA	-	expression tag	UNP P47211
E	332	LEU	-	expression tag	UNP P47211
E	333	GLU	-	expression tag	UNP P47211
E	334	VAL	-	expression tag	UNP P47211
E	335	LEU	-	expression tag	UNP P47211
E	336	PHE	-	expression tag	UNP P47211
E	337	GLN	-	expression tag	UNP P47211

- Molecule 5 is a protein called Guanine nucleotide-binding protein G(I)/G(S)/G(O) subunit gamma-2.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	G	56	Total	C	N	O	S	0	0
			429	269	76	81	3		

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
G	68	SER	CYS	engineered mutation	UNP P59768

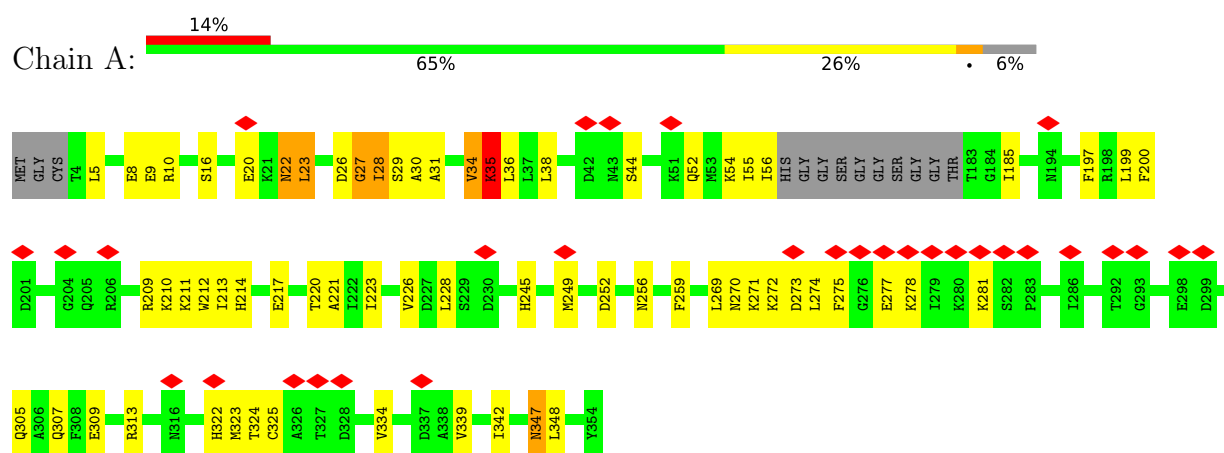
- Molecule 6 is a protein called single Fab chain (svFv16).

Mol	Chain	Residues	Atoms					AltConf	Trace
6	S	231	Total	C	N	O	S	0	0
			1777	1128	294	345	10		

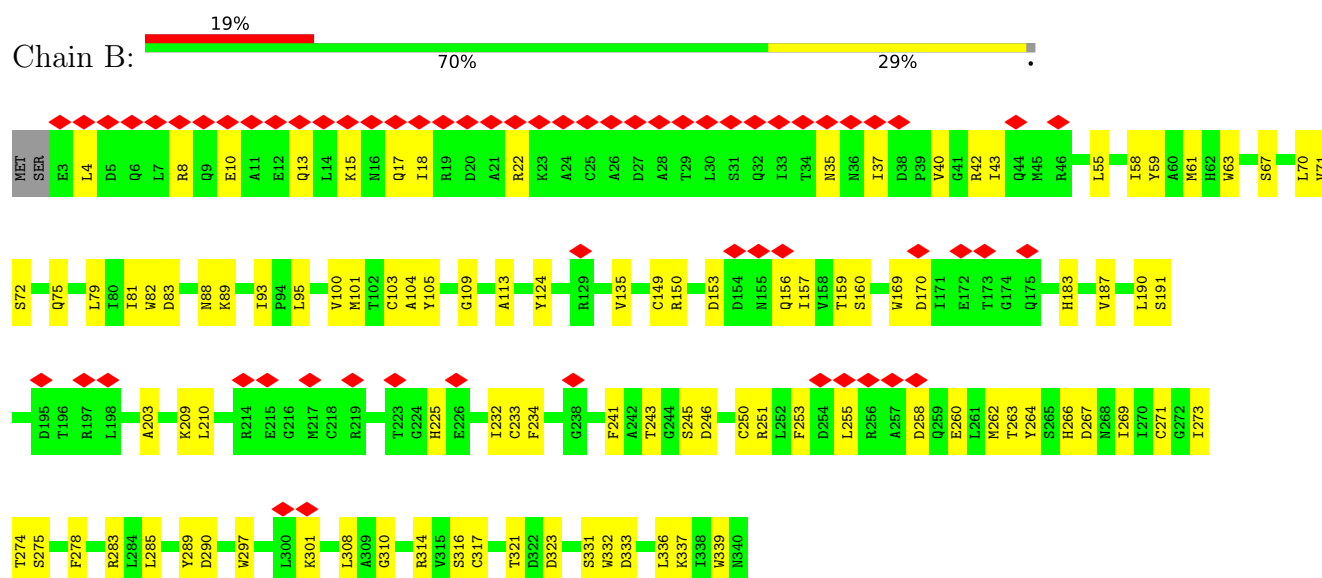
3 Residue-property plots

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.

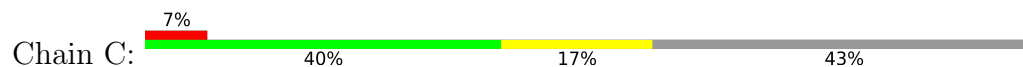
- Molecule 1: G protein subunit alpha o1, Guanine nucleotide-binding protein G(o) subunit alpha



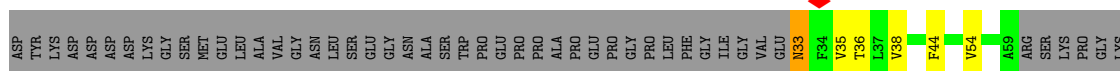
- Molecule 2: Guanine nucleotide-binding protein G(I)/G(S)/G(T) subunit beta-1



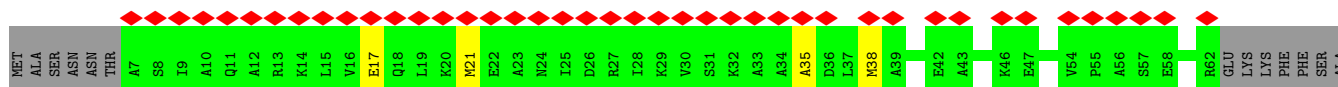
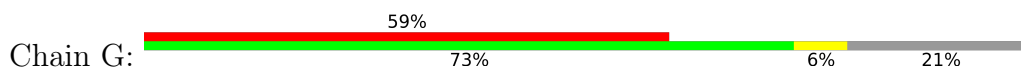
- Molecule 3: Galanin



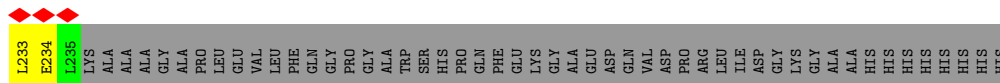
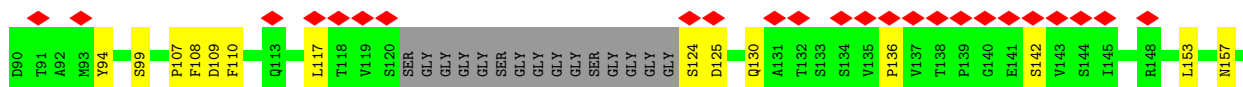
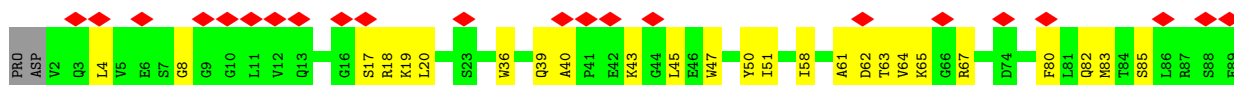
- Molecule 4: Galanin receptor type 1



- Molecule 5: Guanine nucleotide-binding protein G(I)/G(S)/G(O) subunit gamma-2



- Molecule 6: single Fab chain (svFv16)



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	426045	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	50	Depositor
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	2200	Depositor
Magnification	Not provided	
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	3.677	Depositor
Minimum map value	-2.780	Depositor
Average map value	0.001	Depositor
Map value standard deviation	0.102	Depositor
Recommended contour level	0.45	Depositor
Map size (Å)	195.66, 195.66, 195.66	wwPDB
Map dimensions	180, 180, 180	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.087, 1.087, 1.087	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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	0.75	10/1720 (0.6%)	0.70	7/2320 (0.3%)
2	B	0.16	0/2644	0.39	0/3585
3	C	0.11	0/125	0.26	0/171
4	E	0.58	5/2159 (0.2%)	0.68	11/2956 (0.4%)
5	G	0.13	0/435	0.46	0/587
6	S	0.12	0/1821	0.35	0/2469
All	All	0.45	15/8904 (0.2%)	0.54	18/12088 (0.1%)

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	23	LEU	C-O	-9.02	1.13	1.24
1	A	23	LEU	CA-C	-8.76	1.41	1.52
1	A	22	ASN	C-O	-8.48	1.14	1.24
1	A	26	ASP	CA-C	-7.00	1.43	1.52
1	A	26	ASP	C-O	-6.84	1.16	1.24
1	A	34	VAL	C-O	-6.83	1.16	1.24
4	E	164	ILE	CA-CB	-6.69	1.45	1.54
4	E	168	SER	N-CA	-6.44	1.40	1.46
1	A	35	LYS	C-N	-6.08	1.25	1.33
4	E	168	SER	CA-C	-5.67	1.45	1.52
1	A	30	ALA	C-O	-5.54	1.17	1.24
4	E	173	HIS	N-CA	-5.43	1.40	1.46
1	A	31	ALA	C-O	-5.39	1.17	1.24
1	A	31	ALA	CA-C	-5.36	1.45	1.52
4	E	164	ILE	N-CA	-5.01	1.40	1.46

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	35	LYS	N-CA-C	13.38	131.10	110.36
4	E	157	GLY	N-CA-C	-9.12	101.71	112.83
4	E	169	PRO	CA-N-CD	-7.75	101.15	112.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	E	158	CYS	CA-CB-SG	-7.08	98.13	114.40
1	A	35	LYS	CA-C-O	7.05	130.35	121.82
4	E	168	SER	CA-C-N	-6.78	111.53	119.05
4	E	168	SER	C-N-CA	-6.78	111.53	119.05
1	A	23	LEU	CA-C-O	-6.75	113.40	120.55
4	E	166	MET	CB-CA-C	5.79	121.78	110.67
1	A	35	LYS	CA-C-N	-5.78	108.94	121.87
1	A	35	LYS	C-N-CA	-5.78	108.94	121.87
4	E	168	SER	O-C-N	5.54	125.41	120.48
1	A	27	GLY	N-CA-C	-5.52	106.10	112.50
4	E	170	VAL	CB-CA-C	-5.28	105.21	111.97
4	E	164	ILE	N-CA-C	5.26	117.62	111.05
4	E	168	SER	C-N-CD	-5.14	103.92	125.00
1	A	28	ILE	N-CA-C	5.13	115.85	110.62
4	E	166	MET	N-CA-C	5.05	117.68	111.82

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1691	0	1656	50	0
2	B	2597	0	2503	77	0
3	C	121	0	117	6	0
4	E	2095	0	2084	104	0
5	G	429	0	441	2	0
6	S	1777	0	1714	48	0
All	All	8710	0	8515	275	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 16.

All (275) 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)
4:E:113:TYR:CE1	4:E:167:ALA:HB1	1.36	1.60
4:E:113:TYR:CE1	4:E:167:ALA:CB	1.99	1.44
1:A:35:LYS:O	1:A:36:LEU:CG	1.77	1.30
1:A:35:LYS:O	1:A:36:LEU:HG	1.35	1.20
4:E:113:TYR:CD1	4:E:167:ALA:HB1	1.79	1.16
1:A:35:LYS:O	1:A:36:LEU:CD2	1.95	1.14
4:E:168:SER:OG	4:E:169:PRO:HD2	1.50	1.09
4:E:113:TYR:OH	4:E:164:ILE:HD12	1.56	1.06
1:A:35:LYS:O	1:A:36:LEU:HD23	1.61	1.00
4:E:116:THR:OG1	4:E:170:VAL:HG21	1.66	0.95
4:E:113:TYR:CE1	4:E:167:ALA:HB3	2.02	0.95
4:E:117:VAL:CG2	4:E:167:ALA:HB2	2.00	0.92
4:E:113:TYR:HE1	4:E:167:ALA:CB	1.77	0.91
4:E:117:VAL:HG23	4:E:167:ALA:HB2	1.51	0.90
4:E:113:TYR:CE1	4:E:167:ALA:C	2.51	0.88
4:E:113:TYR:HE1	4:E:167:ALA:C	1.82	0.87
4:E:168:SER:OG	4:E:169:PRO:CD	2.23	0.87
4:E:113:TYR:CZ	4:E:167:ALA:HB1	2.12	0.85
4:E:113:TYR:CD1	4:E:167:ALA:O	2.30	0.85
4:E:124:PHE:HB3	4:E:159:ILE:CG2	2.08	0.84
4:E:113:TYR:HD1	4:E:167:ALA:O	1.63	0.82
4:E:166:MET:N	4:E:166:MET:SD	2.52	0.81
6:S:168:ARG:HH12	6:S:174:GLN:HE22	1.30	0.79
4:E:176:LEU:HA	4:E:186:PHE:O	1.83	0.78
1:A:35:LYS:C	1:A:36:LEU:HG	2.08	0.78
1:A:348:LEU:HD11	4:E:227:LEU:HD11	1.64	0.78
2:B:71:VAL:HG12	2:B:81:ILE:HG12	1.63	0.77
4:E:113:TYR:CZ	4:E:167:ALA:CB	2.67	0.76
4:E:120:LEU:HD11	4:E:166:MET:HB2	1.67	0.76
4:E:113:TYR:OH	4:E:164:ILE:CD1	2.33	0.74
2:B:71:VAL:HG13	2:B:105:TYR:HD2	1.53	0.74
2:B:316:SER:OG	2:B:332:TRP:NE1	2.21	0.74
4:E:117:VAL:HG13	4:E:163:SER:OG	1.88	0.74
4:E:120:LEU:CD1	4:E:166:MET:HB2	2.17	0.73
6:S:63:THR:HG23	6:S:64:VAL:HG13	1.70	0.73
6:S:168:ARG:HH21	6:S:210:GLU:HG3	1.53	0.73
6:S:17:SER:OG	6:S:82:GLN:NE2	2.22	0.72
5:G:17:GLU:HG2	5:G:21:MET:HE1	1.73	0.71
1:A:34:VAL:HG11	1:A:339:VAL:HG11	1.72	0.70
4:E:124:PHE:CB	4:E:159:ILE:CG2	2.70	0.70
2:B:271:CYS:HB3	2:B:290:ASP:HB2	1.72	0.70
4:E:117:VAL:HG23	4:E:167:ALA:CB	2.22	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:9:TYR:OH	4:E:112:HIS:NE2	2.26	0.69
2:B:42:ARG:NH1	2:B:43:ILE:O	2.27	0.68
2:B:310:GLY:O	2:B:337:LYS:NZ	2.25	0.68
1:A:252:ASP:OD1	1:A:256:ASN:ND2	2.27	0.67
4:E:113:TYR:CE1	4:E:167:ALA:CA	2.76	0.67
3:C:10:LEU:HD11	4:E:285:ARG:HD3	1.76	0.67
2:B:61:MET:HG3	2:B:317:CYS:HB2	1.78	0.66
6:S:130:GLN:HE22	6:S:228:GLY:HA3	1.60	0.66
2:B:289:TYR:OH	2:B:297:TRP:NE1	2.30	0.65
1:A:8:GLU:OE2	6:S:163:TYR:OH	2.15	0.65
4:E:133:ARG:NH1	4:E:220:TYR:OH	2.31	0.64
4:E:299:ASN:O	4:E:303:TYR:HB2	1.98	0.64
4:E:259:SER:OG	4:E:295:ASN:ND2	2.32	0.63
4:E:105:ALA:HA	4:E:176:LEU:HD13	1.80	0.63
1:A:27:GLY:HA3	2:B:55:LEU:HD13	1.81	0.63
4:E:44:PHE:HB2	4:E:89:ILE:HG21	1.81	0.63
4:E:124:PHE:HB3	4:E:159:ILE:HG21	1.78	0.63
4:E:178:HIS:ND1	4:E:184:GLN:O	2.32	0.63
2:B:58:ILE:HD13	2:B:336:LEU:HD12	1.81	0.62
2:B:232:ILE:HG13	2:B:243:THR:HG22	1.81	0.62
6:S:124:SER:OG	6:S:125:ASP:N	2.31	0.62
2:B:260:GLU:OE2	2:B:263:THR:OG1	2.16	0.62
4:E:33:ASN:O	4:E:36:THR:OG1	2.14	0.61
4:E:117:VAL:HG22	4:E:167:ALA:HB2	1.83	0.61
4:E:155:GLY:O	4:E:159:ILE:HG12	2.00	0.61
4:E:117:VAL:CG2	4:E:167:ALA:CB	2.78	0.61
1:A:211:LYS:O	1:A:214:HIS:NE2	2.34	0.61
1:A:271:LYS:H	1:A:325:CYS:HA	1.65	0.60
2:B:251:ARG:NH1	2:B:263:THR:OG1	2.34	0.60
6:S:39:GLN:HB2	6:S:45:LEU:HD23	1.84	0.60
2:B:156:GLN:NE2	2:B:170:ASP:OD1	2.35	0.60
4:E:169:PRO:HB2	4:E:200:TYR:CZ	2.37	0.60
1:A:272:LYS:HD2	1:A:323:MET:HG2	1.83	0.59
4:E:193:ASP:HB3	4:E:196:HIS:CD2	2.37	0.59
4:E:113:TYR:CD1	4:E:167:ALA:C	2.79	0.59
4:E:124:PHE:CB	4:E:159:ILE:HG22	2.33	0.58
2:B:275:SER:OG	2:B:316:SER:O	2.22	0.58
5:G:35:ALA:HA	5:G:38:MET:SD	2.43	0.58
4:E:121:VAL:HG11	4:E:160:TRP:CZ2	2.38	0.58
2:B:251:ARG:NH1	2:B:260:GLU:OE2	2.34	0.57
1:A:228:LEU:HD11	1:A:269:LEU:HB3	1.85	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:322:HIS:HB2	1:A:334:VAL:HG21	1.86	0.57
6:S:40:ALA:HB3	6:S:43:LYS:HB2	1.86	0.57
6:S:168:ARG:HH12	6:S:174:GLN:NE2	2.01	0.57
6:S:8:GLY:O	6:S:18:ARG:NH1	2.36	0.57
4:E:124:PHE:CB	4:E:159:ILE:HG21	2.35	0.57
4:E:259:SER:HB2	4:E:295:ASN:HB2	1.87	0.57
2:B:75:GLN:NE2	2:B:100:VAL:O	2.38	0.56
1:A:245:HIS:O	1:A:249:MET:HG2	2.05	0.56
2:B:187:VAL:HA	2:B:203:ALA:HA	1.86	0.56
6:S:153:LEU:HA	6:S:221:LEU:HD22	1.86	0.56
6:S:168:ARG:HB2	6:S:171:GLN:HB3	1.88	0.56
6:S:166:LEU:HB2	6:S:176:LEU:HD11	1.87	0.56
2:B:72:SER:HB3	2:B:82:TRP:HE1	1.71	0.56
1:A:10:ARG:NH2	6:S:222:GLU:OE1	2.35	0.56
1:A:27:GLY:HA3	2:B:55:LEU:CD1	2.35	0.56
6:S:19:LYS:HD2	6:S:80:PHE:HB3	1.88	0.55
1:A:35:LYS:HD3	1:A:200:PHE:HE2	1.72	0.55
2:B:105:TYR:HE1	2:B:109:GLY:HA2	1.71	0.55
4:E:113:TYR:OH	4:E:164:ILE:HG23	2.07	0.55
1:A:44:SER:O	1:A:270:ASN:ND2	2.40	0.54
2:B:105:TYR:CE1	2:B:109:GLY:HA2	2.41	0.54
2:B:83:ASP:N	2:B:88:ASN:O	2.37	0.54
2:B:290:ASP:HA	2:B:314:ARG:HG3	1.88	0.54
4:E:116:THR:HG1	4:E:170:VAL:HG21	1.72	0.54
1:A:35:LYS:HD3	1:A:200:PHE:CE2	2.43	0.54
2:B:35:ASN:OD1	2:B:301:LYS:NZ	2.39	0.54
1:A:273:ASP:OD1	1:A:274:LEU:N	2.40	0.53
2:B:153:ASP:OD1	2:B:156:GLN:N	2.40	0.53
4:E:113:TYR:CE1	4:E:167:ALA:O	2.59	0.53
6:S:179:ARG:O	6:S:180:MET:HG2	2.08	0.53
2:B:104:ALA:HB3	2:B:113:ALA:HB3	1.91	0.53
4:E:168:SER:CB	4:E:169:PRO:CD	2.85	0.53
4:E:262:PRO:HG2	4:E:291:LEU:HD12	1.91	0.53
4:E:102:VAL:HG23	4:E:103:LEU:HD12	1.91	0.53
4:E:35:VAL:HA	4:E:38:VAL:HG12	1.90	0.53
2:B:274:THR:OG1	2:B:290:ASP:OD1	2.27	0.52
1:A:271:LYS:HD3	1:A:274:LEU:HD12	1.91	0.52
6:S:36:TRP:CZ3	6:S:94:TYR:HB3	2.44	0.52
4:E:156:VAL:O	4:E:157:GLY:C	2.51	0.52
1:A:36:LEU:HB2	1:A:199:LEU:HD23	1.91	0.52
1:A:217:GLU:HG2	1:A:259:PHE:HB3	1.90	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:E:302:ILE:HG13	4:E:303:TYR:HD1	1.75	0.52
4:E:156:VAL:HA	4:E:159:ILE:CG1	2.40	0.51
4:E:156:VAL:HA	4:E:159:ILE:HG12	1.93	0.51
4:E:162:LEU:O	4:E:163:SER:C	2.51	0.51
4:E:113:TYR:CZ	4:E:167:ALA:HB3	2.41	0.51
2:B:274:THR:HG21	2:B:316:SER:HA	1.92	0.51
2:B:59:TYR:HD2	2:B:101:MET:HB3	1.75	0.51
2:B:157:ILE:HG22	2:B:169:TRP:HB2	1.91	0.51
4:E:130:SER:OG	4:E:216:ILE:HG12	2.10	0.50
2:B:262:MET:HE2	2:B:264:TYR:HE1	1.75	0.50
4:E:54:VAL:HA	4:E:310:PHE:HZ	1.77	0.50
1:A:305:GLN:NE2	1:A:309:GLU:OE2	2.45	0.50
2:B:61:MET:HE2	2:B:70:LEU:HD12	1.92	0.50
1:A:55:ILE:HG23	1:A:56:ILE:H	1.75	0.50
2:B:71:VAL:HG13	2:B:105:TYR:CD2	2.41	0.50
2:B:321:THR:OG1	2:B:323:ASP:OD1	2.30	0.49
4:E:155:GLY:O	4:E:158:CYS:HB3	2.12	0.49
2:B:183:HIS:CE1	2:B:209:LYS:HB2	2.47	0.49
3:C:10:LEU:HD13	4:E:275:PHE:CE2	2.48	0.49
4:E:220:TYR:HA	4:E:223:VAL:HG22	1.95	0.49
1:A:34:VAL:O	1:A:34:VAL:HG12	2.12	0.49
2:B:267:ASP:N	2:B:267:ASP:OD1	2.45	0.49
2:B:245:SER:OG	2:B:246:ASP:N	2.46	0.49
1:A:5:LEU:O	1:A:10:ARG:NH1	2.45	0.49
1:A:9:GLU:OE2	6:S:157:ASN:ND2	2.37	0.49
2:B:70:LEU:HD23	2:B:82:TRP:HB2	1.96	0.48
1:A:220:THR:HG21	1:A:339:VAL:HG13	1.95	0.48
2:B:253:PHE:HD1	2:B:260:GLU:HA	1.78	0.48
2:B:250:CYS:SG	2:B:273:ILE:HG12	2.53	0.48
4:E:117:VAL:HG13	4:E:163:SER:HG	1.75	0.48
6:S:4:LEU:HD12	6:S:110:PHE:HD2	1.79	0.48
2:B:10:GLU:O	2:B:13:GLN:HG3	2.14	0.47
4:E:262:PRO:HA	4:E:265:ILE:HG22	1.96	0.47
2:B:234:PHE:CD1	2:B:241:PHE:HB3	2.49	0.47
3:C:2:TRP:HE1	4:E:281:SER:HB3	1.79	0.47
4:E:113:TYR:HB2	4:E:171:ALA:HB2	1.95	0.47
6:S:162:LEU:HD11	6:S:217:CYS:HB2	1.96	0.47
2:B:225:HIS:HE2	2:B:243:THR:HG1	1.59	0.47
1:A:221:ALA:HB2	1:A:342:ILE:HD13	1.95	0.47
2:B:266:HIS:HB3	2:B:269:ILE:HG12	1.97	0.47
2:B:83:ASP:HB3	2:B:88:ASN:HB2	1.96	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:S:61:ALA:HB3	6:S:64:VAL:HG22	1.96	0.46
2:B:191:SER:HB3	2:B:232:ILE:HG23	1.97	0.46
4:E:159:ILE:HG12	4:E:159:ILE:H	1.46	0.46
6:S:51:ILE:HG13	6:S:58:ILE:HG13	1.96	0.46
6:S:232:LYS:HD2	6:S:233:LEU:N	2.30	0.46
2:B:63:TRP:CD2	2:B:321:THR:HG22	2.51	0.46
6:S:166:LEU:HD13	6:S:215:TYR:CE1	2.50	0.46
6:S:99:SER:OG	6:S:108:PHE:N	2.43	0.46
6:S:168:ARG:NE	6:S:210:GLU:OE2	2.49	0.46
4:E:148:VAL:HG12	4:E:151:ASN:H	1.81	0.46
4:E:164:ILE:HD13	4:E:164:ILE:HA	1.56	0.46
4:E:119:MET:O	4:E:123:ILE:HG13	2.15	0.46
4:E:168:SER:HG	4:E:169:PRO:HD2	1.75	0.46
1:A:16:SER:O	1:A:20:GLU:HG2	2.15	0.46
2:B:22:ARG:NH2	2:B:258:ASP:OD1	2.49	0.46
2:B:61:MET:HE1	2:B:63:TRP:CD1	2.51	0.46
6:S:162:LEU:HD22	6:S:200:PHE:CG	2.51	0.45
6:S:67:ARG:NH1	6:S:85:SER:O	2.50	0.45
1:A:313:ARG:HG2	1:A:313:ARG:O	2.16	0.45
6:S:47:TRP:HZ2	6:S:50:TYR:HB3	1.82	0.45
4:E:234:MET:HE2	4:E:234:MET:HA	1.99	0.45
6:S:20:LEU:HG	6:S:83:MET:HE1	1.98	0.45
6:S:62:ASP:HA	6:S:65:LYS:NZ	2.32	0.45
2:B:67:SER:O	2:B:67:SER:OG	2.31	0.44
4:E:164:ILE:HD12	4:E:164:ILE:HG23	1.52	0.44
6:S:19:LYS:HE2	6:S:82:GLN:OE1	2.17	0.44
6:S:164:TRP:CZ3	6:S:217:CYS:HB3	2.52	0.44
6:S:142:SER:HA	6:S:204:ILE:O	2.17	0.44
6:S:160:THR:HB	6:S:180:MET:HE2	1.98	0.44
4:E:177:PHE:HD1	4:E:190:GLN:HE22	1.64	0.44
2:B:233:CYS:SG	2:B:234:PHE:N	2.90	0.44
1:A:22:ASN:N	1:A:22:ASN:HD22	2.16	0.44
4:E:270:ALA:HB2	4:E:275:PHE:CE2	2.53	0.44
2:B:331:SER:OG	2:B:332:TRP:N	2.51	0.44
4:E:283:LEU:HA	4:E:286:ILE:HG22	2.00	0.44
2:B:79:LEU:HD23	2:B:93:ILE:HD12	1.99	0.43
4:E:178:HIS:CE1	4:E:185:THR:HB	2.52	0.43
6:S:62:ASP:HA	6:S:65:LYS:HZ1	1.83	0.43
4:E:162:LEU:HD23	4:E:162:LEU:HA	1.64	0.43
1:A:52:GLN:HB3	1:A:54:LYS:NZ	2.33	0.43
2:B:40:VAL:HG23	2:B:283:ARG:HH21	1.82	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:E:113:TYR:CD1	4:E:167:ALA:CB	2.66	0.43
1:A:209:ARG:O	1:A:213:ILE:HB	2.18	0.43
2:B:187:VAL:HG12	2:B:203:ALA:HB2	1.99	0.43
4:E:270:ALA:HB2	4:E:275:PHE:HE2	1.83	0.43
1:A:275:PHE:HA	1:A:278:LYS:HG2	2.00	0.43
1:A:185:ILE:HD11	1:A:200:PHE:HB3	2.01	0.43
2:B:149:CYS:HB3	2:B:159:THR:HG22	2.00	0.43
4:E:261:LEU:HD12	4:E:261:LEU:HA	1.90	0.43
1:A:20:GLU:OE2	2:B:89:LYS:NZ	2.26	0.43
1:A:277:GLU:O	1:A:281:LYS:HG2	2.19	0.43
2:B:124:TYR:CE2	2:B:135:VAL:HG22	2.53	0.43
6:S:108:PHE:CZ	6:S:218:MET:HE1	2.54	0.42
2:B:308:LEU:HD23	2:B:339:TRP:CD2	2.55	0.42
4:E:166:MET:O	4:E:200:TYR:OH	2.35	0.42
2:B:210:LEU:HD11	2:B:255:LEU:HD12	2.00	0.42
4:E:156:VAL:C	4:E:158:CYS:N	2.75	0.42
6:S:109:ASP:OD1	6:S:109:ASP:N	2.45	0.42
1:A:226:VAL:HB	1:A:269:LEU:HD23	2.01	0.42
1:A:347:ASN:HD22	1:A:347:ASN:HA	1.63	0.42
2:B:15:LYS:O	2:B:18:ILE:HG22	2.20	0.42
1:A:210:LYS:HB2	1:A:210:LYS:HE3	1.87	0.42
4:E:121:VAL:HG23	4:E:163:SER:HB2	2.01	0.42
6:S:162:LEU:HD22	6:S:200:PHE:CD2	2.55	0.42
3:C:11:LEU:HB3	3:C:15:ALA:HB3	2.01	0.42
4:E:90:PRO:O	4:E:94:THR:HG23	2.19	0.42
4:E:211:LEU:HB2	4:E:212:PRO:HD3	2.02	0.42
4:E:236:LYS:O	4:E:239:GLU:HG3	2.20	0.42
6:S:107:PRO:O	6:S:175:LEU:HD22	2.20	0.42
2:B:95:LEU:HD13	2:B:100:VAL:HG11	2.02	0.42
4:E:113:TYR:HE1	4:E:167:ALA:CA	2.22	0.42
1:A:35:LYS:HB3	1:A:36:LEU:H	1.70	0.42
6:S:220:HIS:HA	6:S:225:LEU:HD22	2.02	0.42
2:B:103:CYS:O	2:B:150:ARG:NH1	2.53	0.41
2:B:160:SER:OG	2:B:190:LEU:HD23	2.20	0.41
4:E:148:VAL:HB	4:E:151:ASN:HB2	2.01	0.41
1:A:272:LYS:HZ2	1:A:324:THR:C	2.27	0.41
2:B:63:TRP:CE2	2:B:321:THR:HG22	2.54	0.41
6:S:136:PRO:HA	6:S:234:GLU:HG2	2.01	0.41
6:S:177:ILE:HG13	6:S:183:LEU:HD12	2.03	0.41
4:E:206:VAL:O	4:E:210:LEU:HB2	2.20	0.41
6:S:219:GLN:OE1	6:S:221:LEU:N	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:4:LEU:O	2:B:8:ARG:HG3	2.19	0.41
3:C:10:LEU:HD11	4:E:285:ARG:CD	2.48	0.41
4:E:164:ILE:O	4:E:164:ILE:HG22	2.19	0.41
4:E:266:ILE:HD12	4:E:285:ARG:HD2	2.02	0.41
2:B:160:SER:HB3	2:B:187:VAL:HG23	2.01	0.41
2:B:160:SER:HB3	2:B:187:VAL:CG2	2.50	0.41
4:E:124:PHE:HB2	4:E:159:ILE:CG2	2.48	0.41
4:E:123:ILE:HG22	4:E:212:PRO:HG2	2.03	0.41
1:A:38:LEU:HD23	1:A:223:ILE:HB	2.03	0.41
2:B:13:GLN:O	2:B:17:GLN:OE1	2.39	0.41
2:B:37:ILE:HG21	2:B:283:ARG:HH12	1.86	0.41
2:B:225:HIS:NE2	2:B:243:THR:OG1	2.50	0.41
2:B:278:PHE:CD1	2:B:285:LEU:HD13	2.56	0.41
4:E:123:ILE:HG21	4:E:208:GLY:O	2.20	0.41
4:E:262:PRO:O	4:E:265:ILE:HG22	2.20	0.41
6:S:18:ARG:HG2	6:S:117:LEU:HD11	2.03	0.41
6:S:215:TYR:O	6:S:230:GLY:HA2	2.21	0.41
2:B:266:HIS:HB2	2:B:297:TRP:HH2	1.85	0.41
2:B:314:ARG:C	2:B:331:SER:HG	2.24	0.41
4:E:84:TYR:CE2	4:E:119:MET:HB2	2.55	0.41
6:S:163:TYR:HE1	6:S:220:HIS:HB2	1.86	0.41
1:A:197:PHE:HE1	1:A:339:VAL:HG21	1.85	0.40
1:A:23:LEU:N	1:A:23:LEU:HD23	2.34	0.40
2:B:331:SER:HG	2:B:332:TRP:H	1.69	0.40
2:B:333:ASP:N	2:B:333:ASP:OD1	2.52	0.40
1:A:209:ARG:HG2	1:A:212:TRP:CZ2	2.56	0.40
4:E:126:LEU:HD22	4:E:303:TYR:CZ	2.56	0.40
4:E:76:ASN:C	4:E:76:ASN:HD22	2.27	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	211/228 (92%)	196 (93%)	15 (7%)	0	100	100
2	B	336/340 (99%)	311 (93%)	25 (7%)	0	100	100
3	C	15/30 (50%)	13 (87%)	2 (13%)	0	100	100
4	E	261/347 (75%)	255 (98%)	6 (2%)	0	100	100
5	G	54/71 (76%)	51 (94%)	3 (6%)	0	100	100
6	S	227/298 (76%)	222 (98%)	5 (2%)	0	100	100
All	All	1104/1314 (84%)	1048 (95%)	56 (5%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

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

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	184/195 (94%)	179 (97%)	5 (3%)	39	63
2	B	280/283 (99%)	280 (100%)	0	100	100
3	C	11/23 (48%)	11 (100%)	0	100	100
4	E	223/292 (76%)	219 (98%)	4 (2%)	51	70
5	G	45/58 (78%)	45 (100%)	0	100	100
6	S	196/237 (83%)	196 (100%)	0	100	100
All	All	939/1088 (86%)	930 (99%)	9 (1%)	65	76

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

Mol	Chain	Res	Type
1	A	28	ILE
1	A	29	SER
1	A	35	LYS
1	A	307	GLN
1	A	347	ASN
4	E	33	ASN
4	E	159	ILE

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Mol	Chain	Res	Type
4	E	166	MET
4	E	169	PRO

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	22	ASN
1	A	347	ASN
2	B	183	HIS
4	E	92	GLN
4	E	196	HIS
4	E	299	ASN
5	G	44	HIS
6	S	82	GLN
6	S	174	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

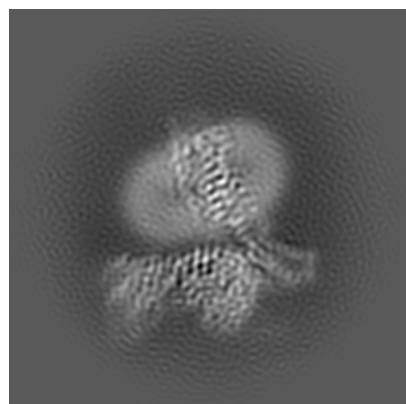
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-33229. These allow visual inspection of the internal detail of the map and identification of artifacts.

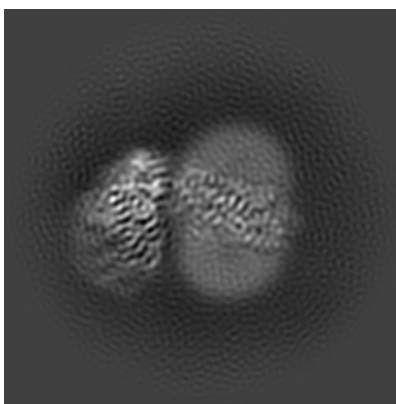
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

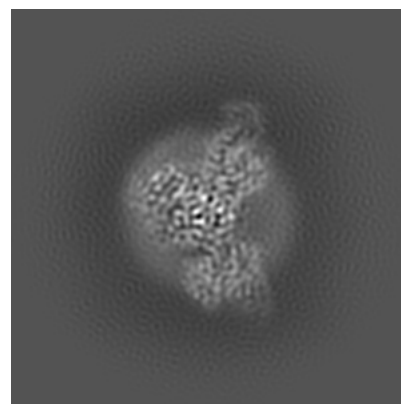
6.1.1 Primary map



X

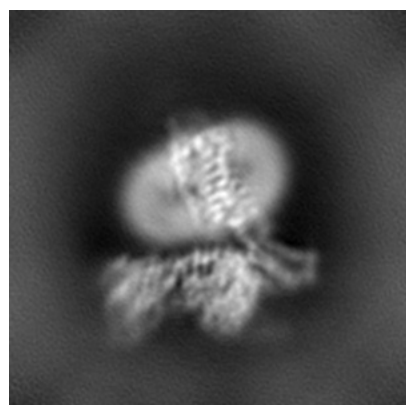


Y

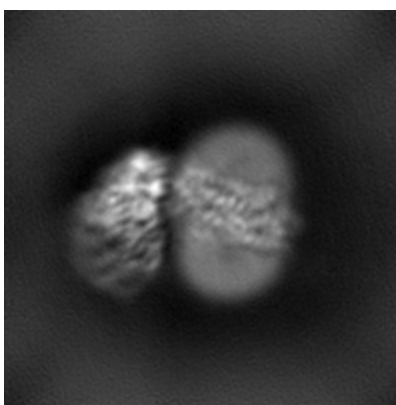


Z

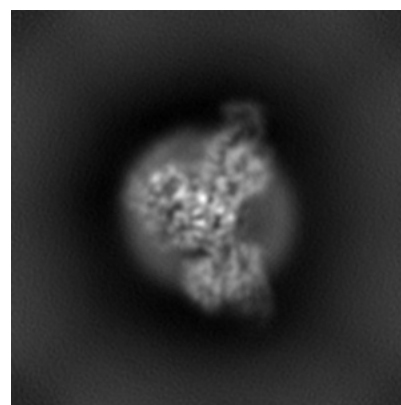
6.1.2 Raw map



X



Y

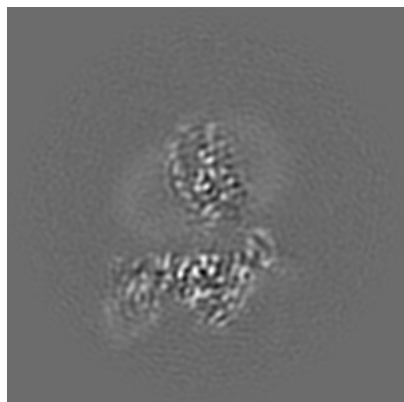


Z

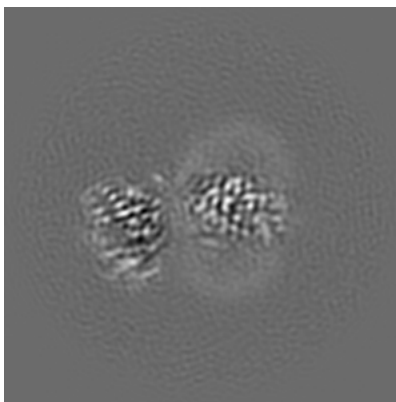
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

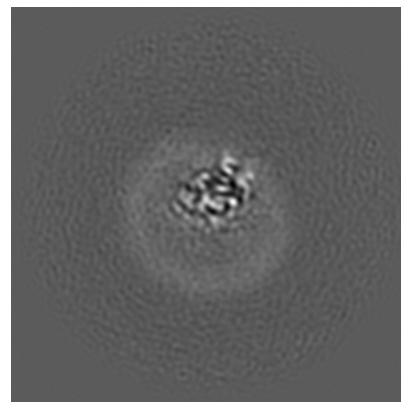
6.2.1 Primary map



X Index: 90

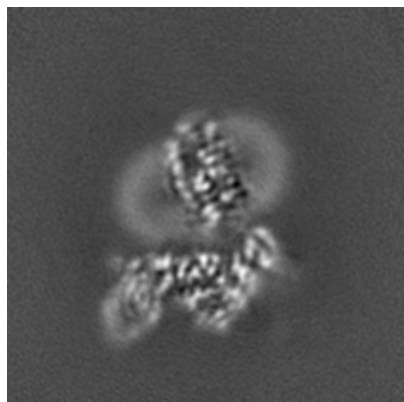


Y Index: 90

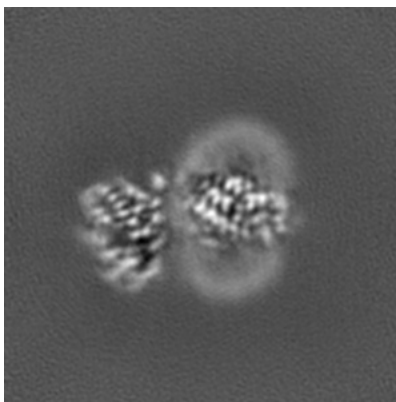


Z Index: 90

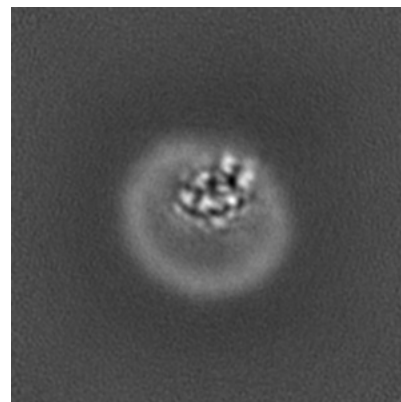
6.2.2 Raw map



X Index: 90



Y Index: 90

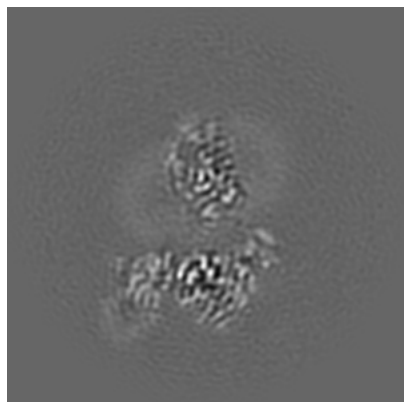


Z Index: 90

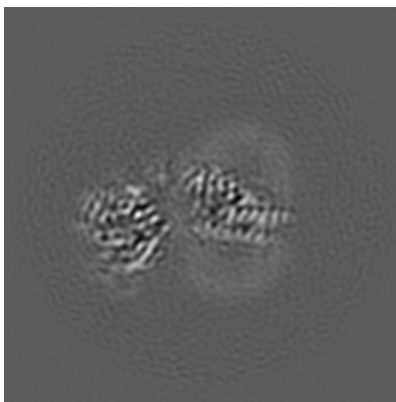
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

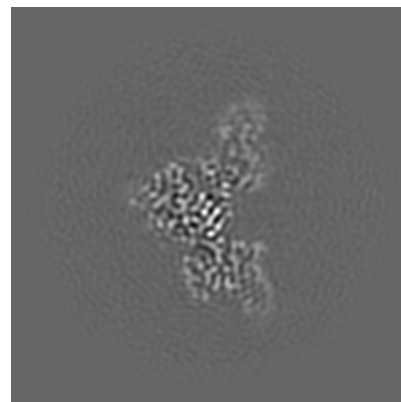
6.3.1 Primary map



X Index: 89

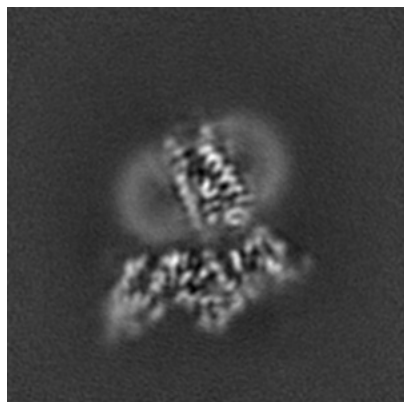


Y Index: 95

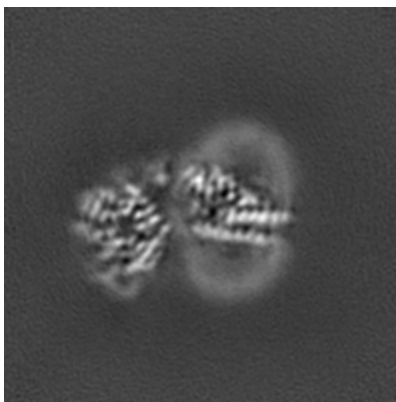


Z Index: 59

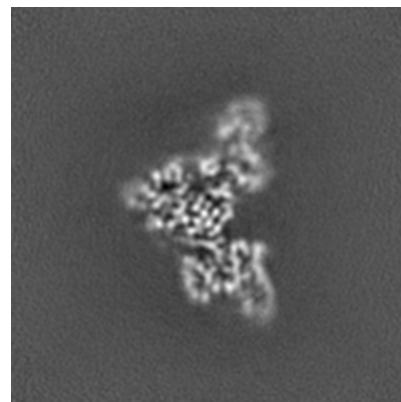
6.3.2 Raw map



X Index: 93



Y Index: 95

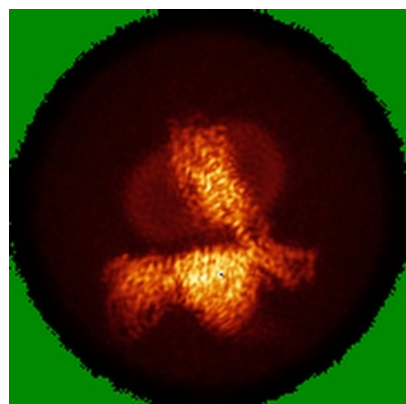


Z Index: 58

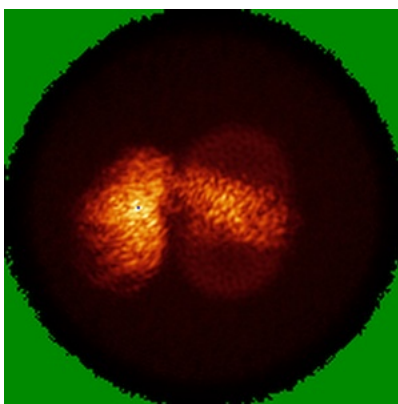
The images above show the largest variance slices of the map in three orthogonal directions.

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

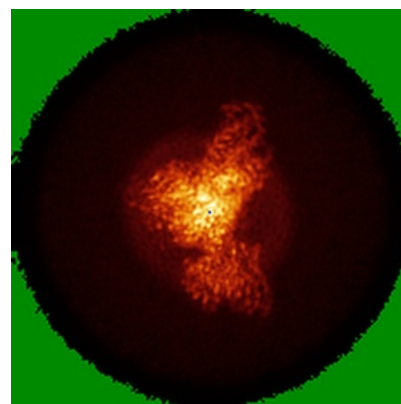
6.4.1 Primary map



X

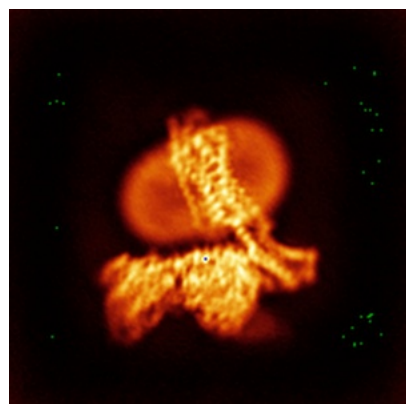


Y

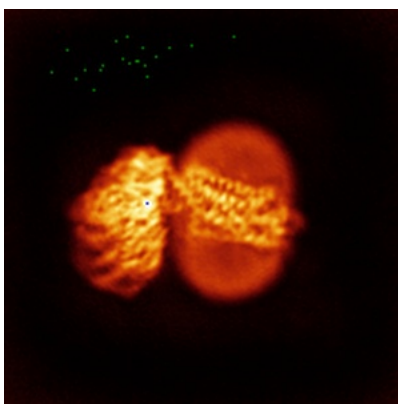


Z

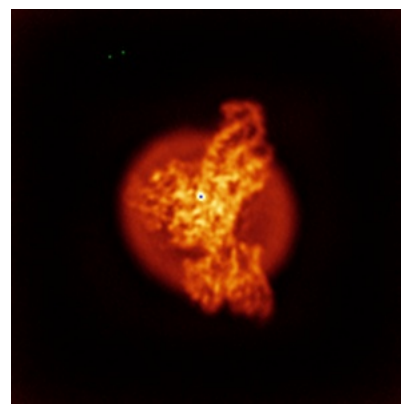
6.4.2 Raw map



X



Y

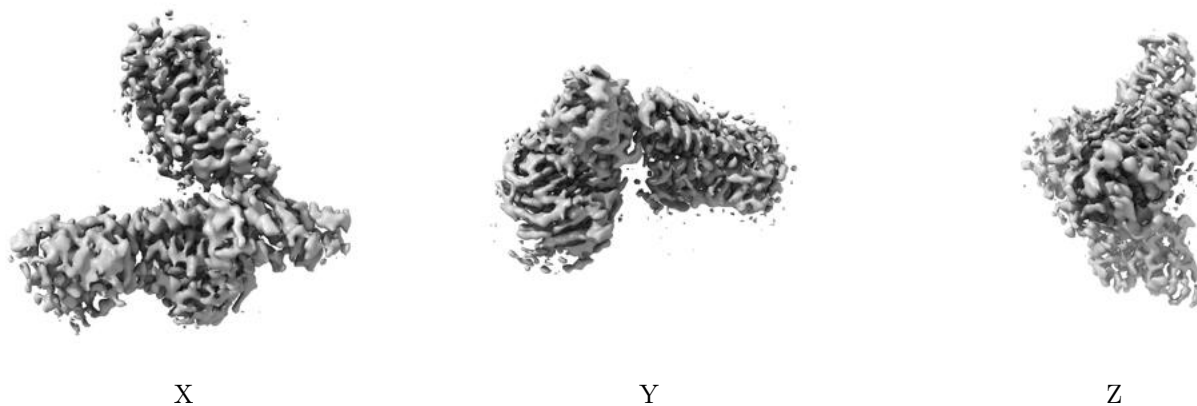


Z

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.

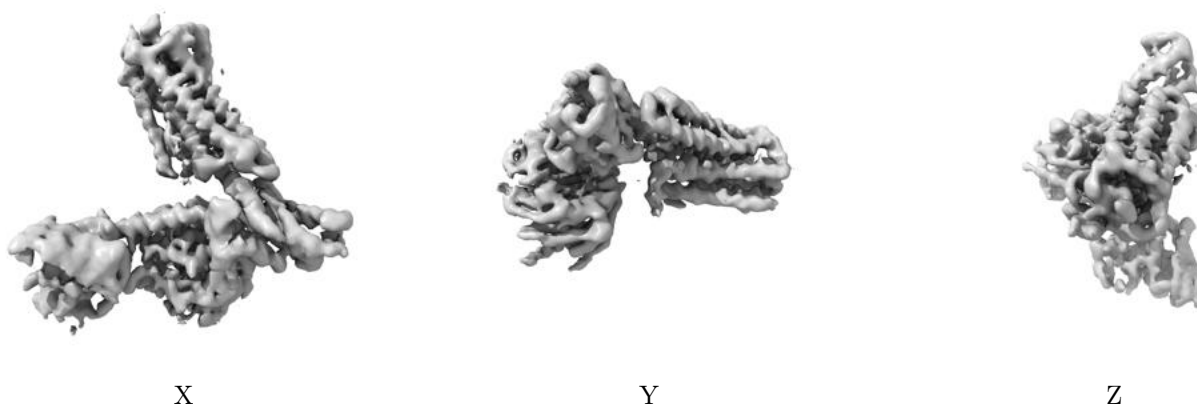
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.45. 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.

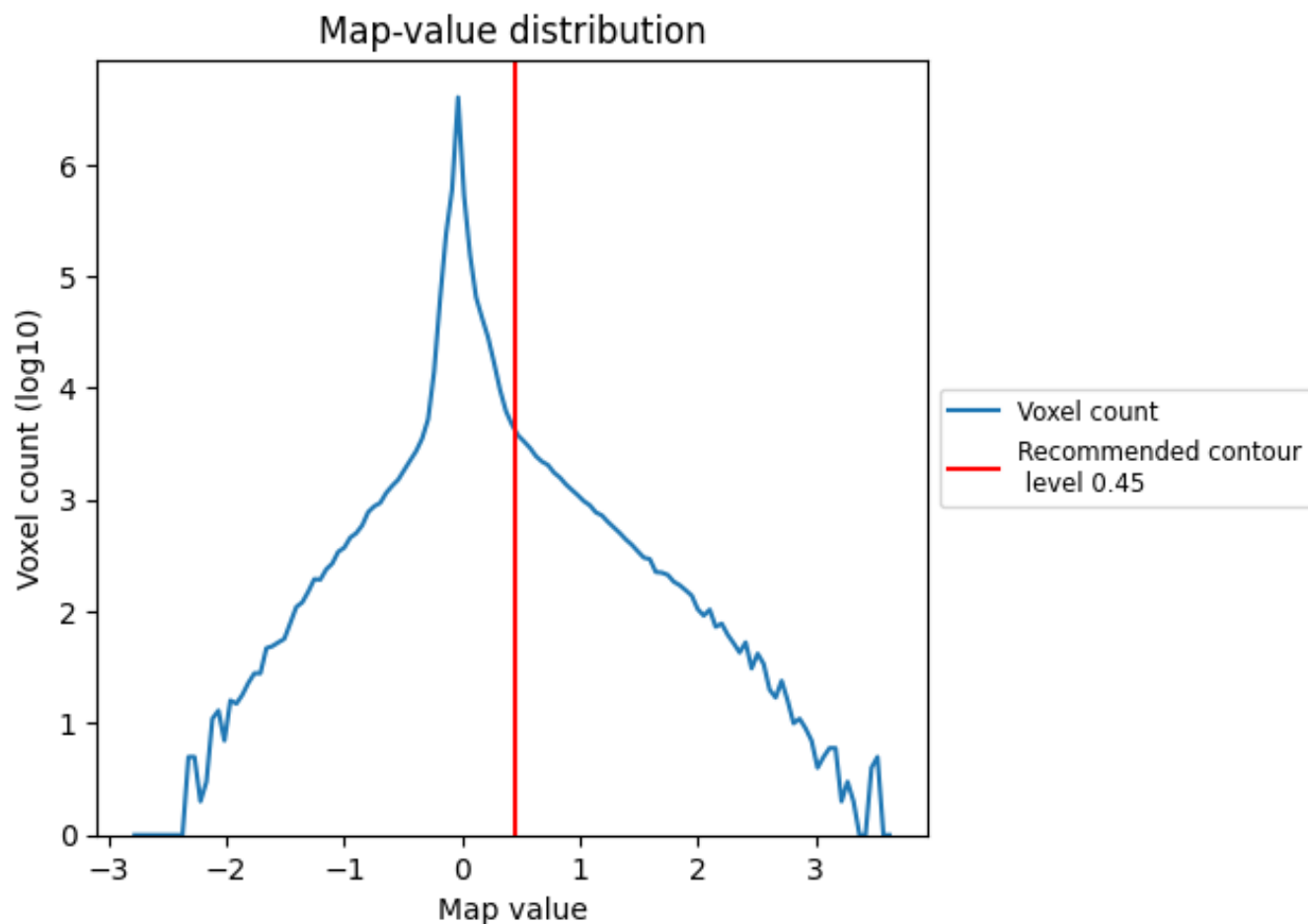
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

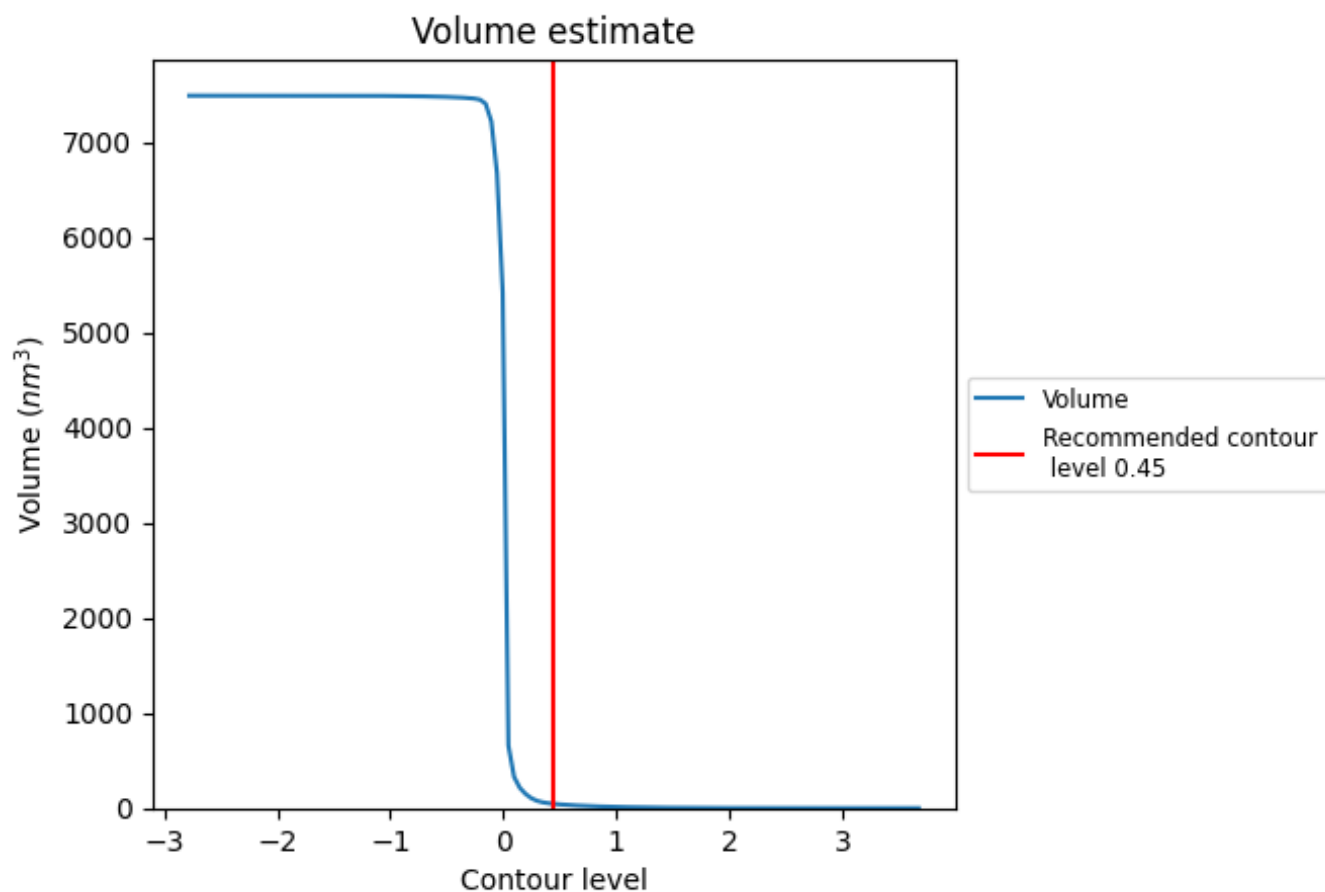
This section contains the results of statistical analysis of the map.

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.

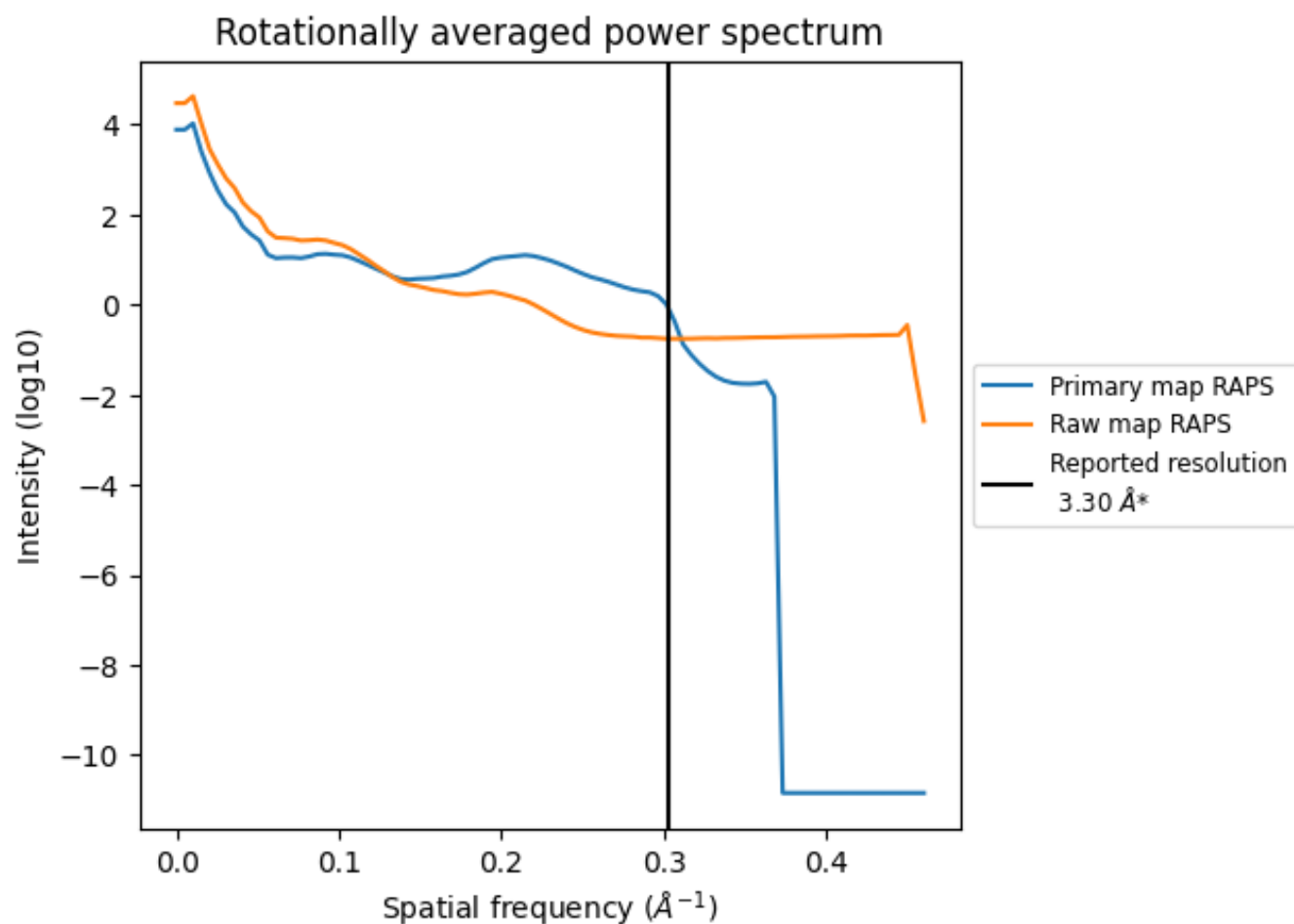
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 45 nm³; this corresponds to an approximate mass of 41 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 ⓘ

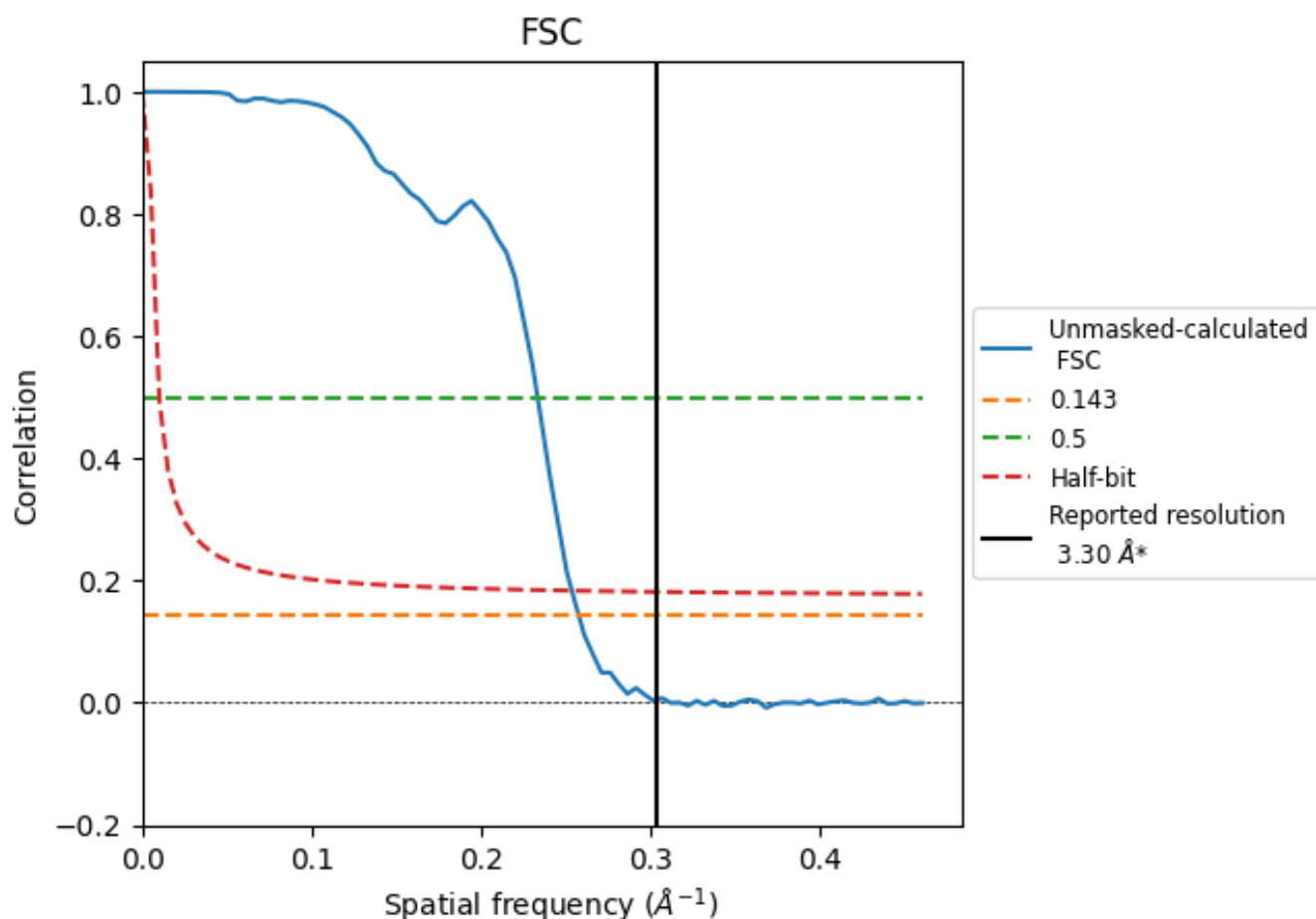


*Reported resolution corresponds to spatial frequency of $0.303 \text{ \AA}^{-1}$

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 0.303 Å⁻¹

8.2 Resolution estimates [i](#)

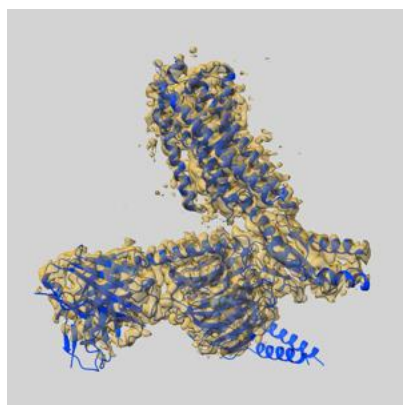
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.30	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	3.88	4.29	3.95

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 3.88 differs from the reported value 3.3 by more than 10 %

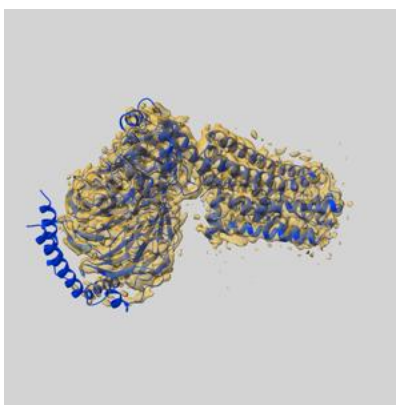
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-33229 and PDB model 7XJJ. Per-residue inclusion information can be found in section [3](#) on page [7](#).

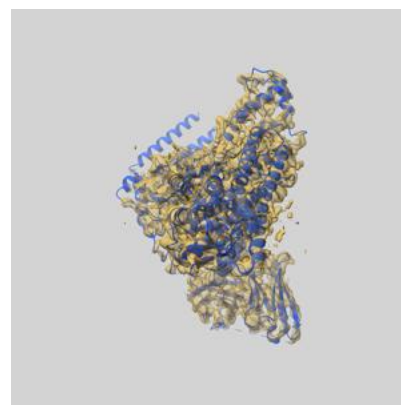
9.1 Map-model overlay [i](#)



X



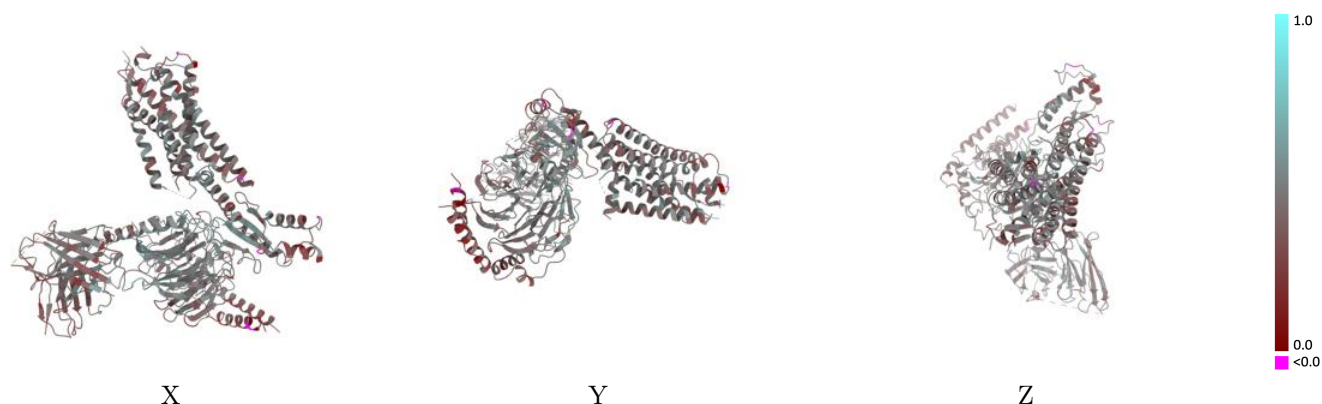
Y



Z

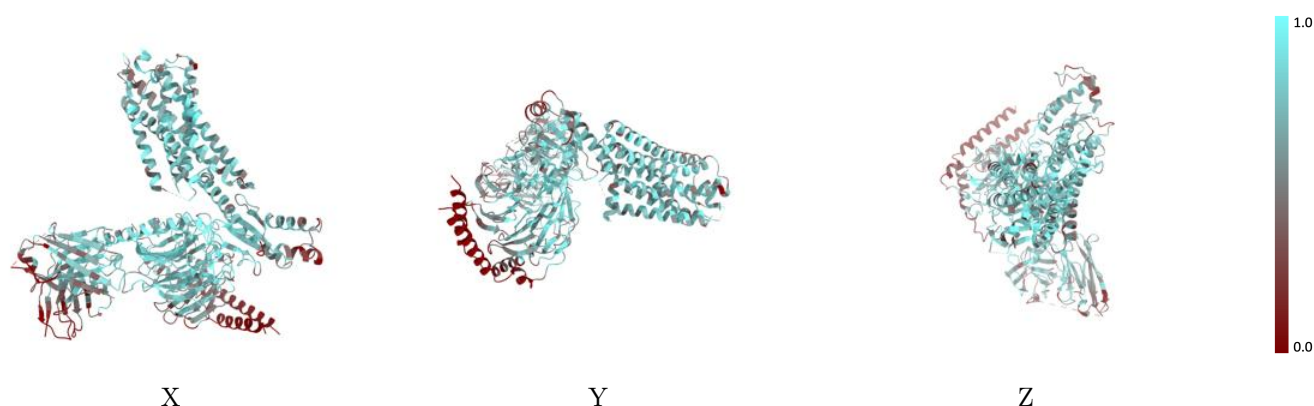
The images above show the 3D surface view of the map at the recommended contour level 0.45 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



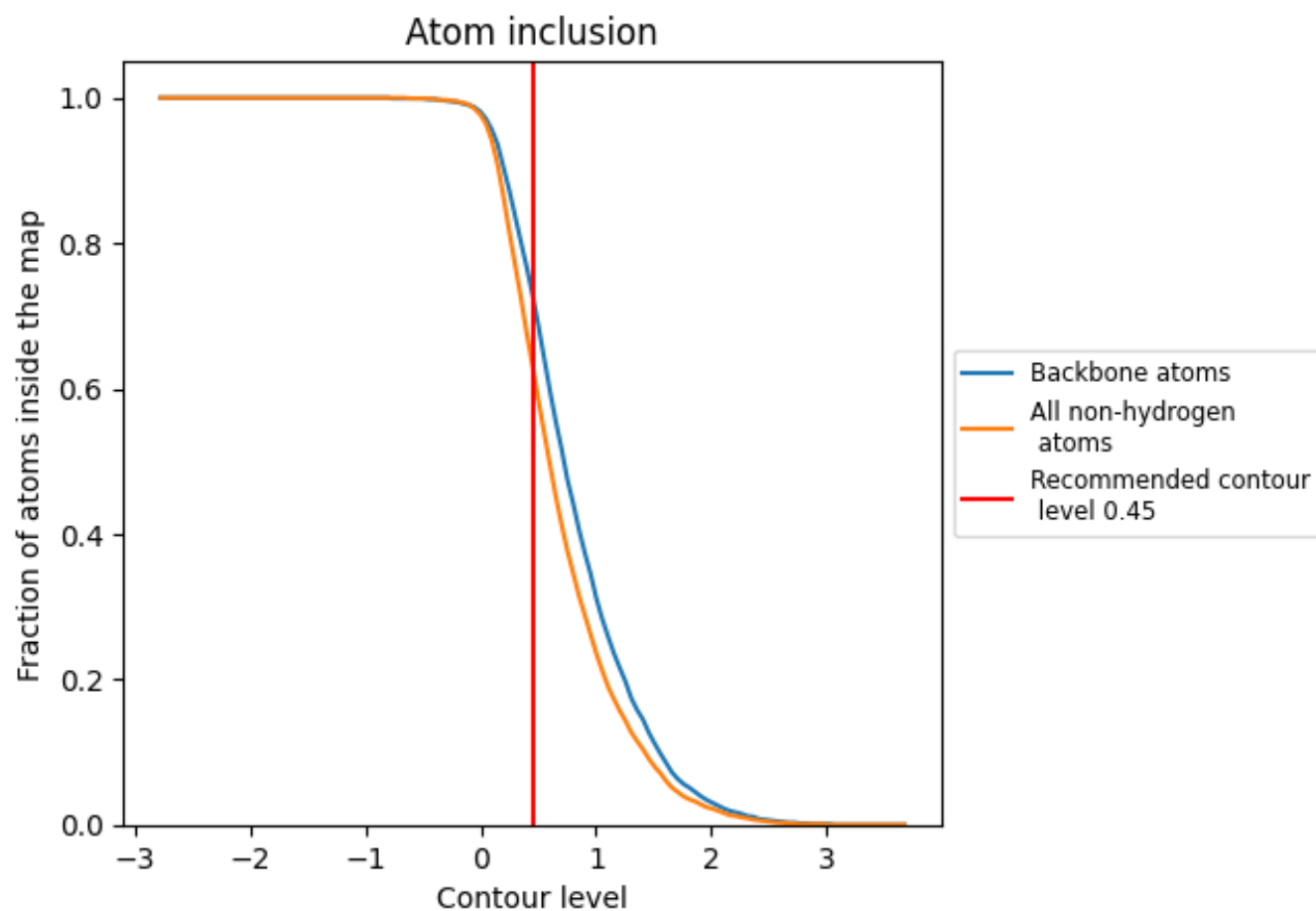
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.45).

9.4 Atom inclusion [i](#)



At the recommended contour level, 73% of all backbone atoms, 63% 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 (0.45) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.6300	0.4360
A	0.6730	0.4530
B	0.6410	0.4480
C	0.7400	0.4380
E	0.7060	0.4300
G	0.2540	0.3310
S	0.5640	0.4320

1.0

0.0

<0.0