



Full wwPDB EM Validation Report ⓘ

Oct 14, 2024 – 07:04 PM JST

PDB ID : 7W01
EMDB ID : EMD-32233
Title : Cryo-EM structure of nucleotide-free ABCA3
Authors : Xie, T.; Zhang, Z.K.; Yue, J.; Gong, X.
Deposited on : 2021-11-17
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.dev113
Mogul : 1.8.5 (274361), CSD as541be (2020)
MolProbity : 4.02b-467
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) : 2.39

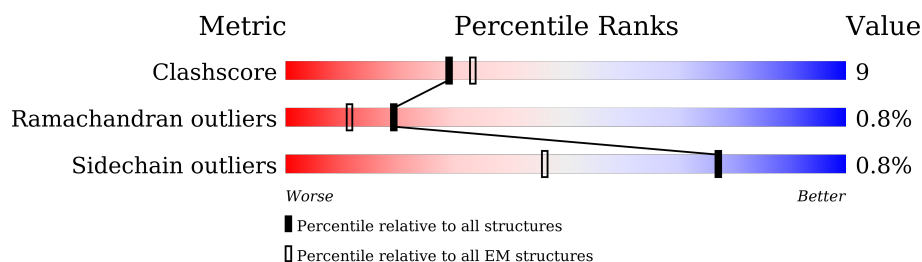
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY


The reported resolution of this entry is 3.30 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)
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 $\leq 5\%$. 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	1748	 74% 16% • 9%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	PX4	A	1804	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 12724 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 Phospholipid-transporting ATPase ABCA3.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	1598	Total	C	N	O	S	0	0
			12556	8151	2110	2225	70		

There are 44 discrepancies between the modelled and reference sequences:

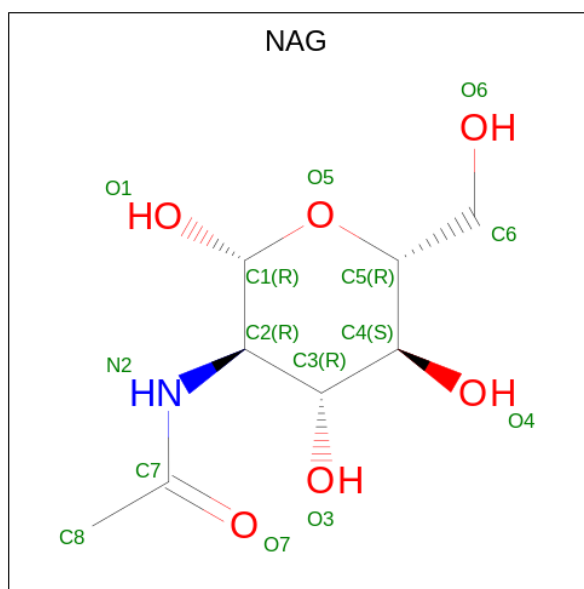
Chain	Residue	Modelled	Actual	Comment	Reference
A	-20	MET	-	initiating methionine	UNP Q99758
A	-19	ALA	-	expression tag	UNP Q99758
A	-18	ASP	-	expression tag	UNP Q99758
A	-17	TYR	-	expression tag	UNP Q99758
A	-16	LYS	-	expression tag	UNP Q99758
A	-15	ASP	-	expression tag	UNP Q99758
A	-14	ASP	-	expression tag	UNP Q99758
A	-13	ASP	-	expression tag	UNP Q99758
A	-12	ASP	-	expression tag	UNP Q99758
A	-11	LYS	-	expression tag	UNP Q99758
A	-10	SER	-	expression tag	UNP Q99758
A	-9	GLY	-	expression tag	UNP Q99758
A	-8	PRO	-	expression tag	UNP Q99758
A	-7	ASP	-	expression tag	UNP Q99758
A	-6	GLU	-	expression tag	UNP Q99758
A	-5	VAL	-	expression tag	UNP Q99758
A	-4	ASP	-	expression tag	UNP Q99758
A	-3	ALA	-	expression tag	UNP Q99758
A	-2	SER	-	expression tag	UNP Q99758
A	-1	GLY	-	expression tag	UNP Q99758
A	0	ARG	-	expression tag	UNP Q99758
A	1705	LEU	-	expression tag	UNP Q99758
A	1706	GLU	-	expression tag	UNP Q99758
A	1707	GLY	-	expression tag	UNP Q99758
A	1708	SER	-	expression tag	UNP Q99758
A	1709	ASP	-	expression tag	UNP Q99758
A	1710	GLU	-	expression tag	UNP Q99758
A	1711	VAL	-	expression tag	UNP Q99758

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Chain	Residue	Modelled	Actual	Comment	Reference
A	1712	ASP	-	expression tag	UNP Q99758
A	1713	ALA	-	expression tag	UNP Q99758
A	1714	VAL	-	expression tag	UNP Q99758
A	1715	GLU	-	expression tag	UNP Q99758
A	1716	GLY	-	expression tag	UNP Q99758
A	1717	SER	-	expression tag	UNP Q99758
A	1718	HIS	-	expression tag	UNP Q99758
A	1719	HIS	-	expression tag	UNP Q99758
A	1720	HIS	-	expression tag	UNP Q99758
A	1721	HIS	-	expression tag	UNP Q99758
A	1722	HIS	-	expression tag	UNP Q99758
A	1723	HIS	-	expression tag	UNP Q99758
A	1724	HIS	-	expression tag	UNP Q99758
A	1725	HIS	-	expression tag	UNP Q99758
A	1726	HIS	-	expression tag	UNP Q99758
A	1727	HIS	-	expression tag	UNP Q99758

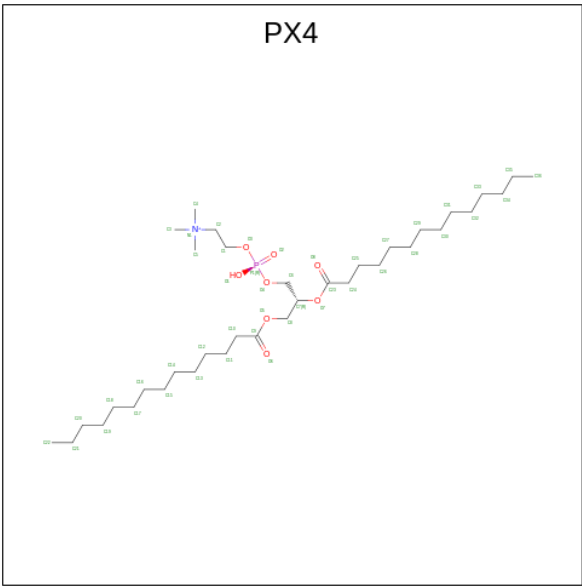
- Molecule 2 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



Mol	Chain	Residues	Atoms				AltConf
2	A	1	Total	C	N	O	0
			14	8	1	5	
2	A	1	Total	C	N	O	0
			14	8	1	5	

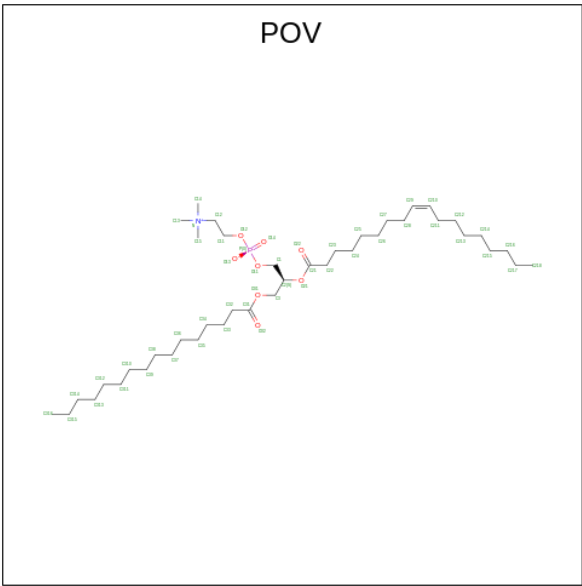
- Molecule 3 is 1,2-DIMYRISTOYL-SN-GLYCERO-3-PHOSPHOCHOLINE (three-letter

code: PX4) (formula: C₃₆H₇₃NO₈P) (labeled as "Ligand of Interest" by depositor).

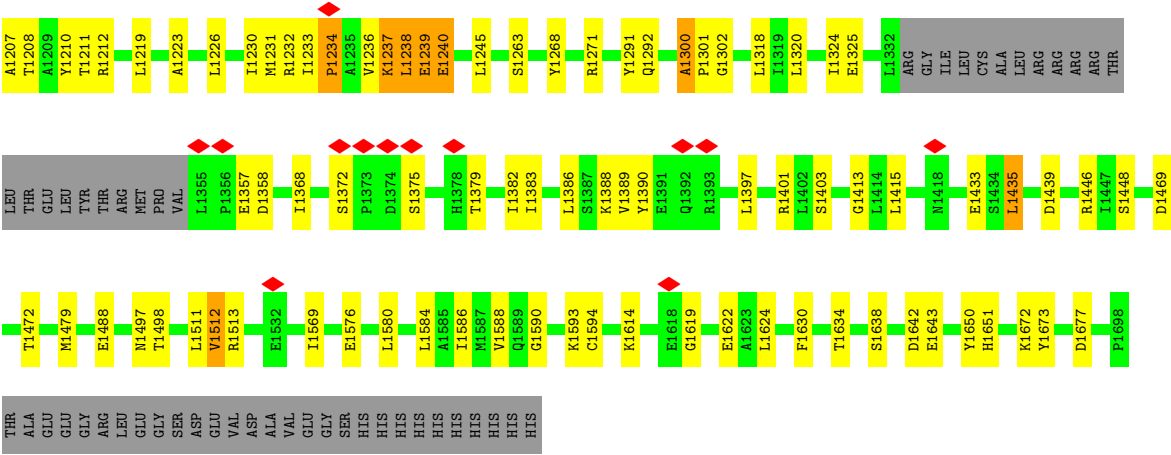


Mol	Chain	Residues	Atoms					AltConf
3	A	1	Total	C	N	O	P	0
			46	36	1	8	1	
3	A	1	Total	C	N	O	P	0
			46	36	1	8	1	

- Molecule 4 is (2S)-3-(hexadecanoyloxy)-2-[(9Z)-octadec-9-enoyloxy]propyl 2-(trimethylamm onio)ethyl phosphate (three-letter code: POV) (formula: C₄₂H₈₂NO₈P) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms		AltConf
4	A	1	Total 16	C 16	0
4	A	1	Total 16	C 16	0
4	A	1	Total 16	C 16	0



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	111435	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	NONE	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)	2000	Depositor
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	0.195	Depositor
Minimum map value	-0.102	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	0.006	Depositor
Recommended contour level	0.025	Depositor
Map size (\AA)	259.2, 259.2, 259.2	wwPDB
Map dimensions	240, 240, 240	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (\AA)	1.08, 1.08, 1.08	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: POV, NAG, PX4

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	0.51	0/12863	0.58	3/17442 (0.0%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	5

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	729	LEU	CA-CB-CG	6.41	130.05	115.30
1	A	955	MET	CA-CB-CG	5.92	123.36	113.30
1	A	1435	LEU	CB-CG-CD1	-5.68	101.35	111.00

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	1300	ALA	Peptide
1	A	140	ASN	Peptide
1	A	1401	ARG	Peptide
1	A	155	ARG	Peptide
1	A	156	PHE	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	12556	0	12608	228	0
2	A	28	0	26	0	0
3	A	92	0	144	29	0
4	A	48	0	78	5	0
All	All	12724	0	12856	233	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 (233) 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:370:LYS:HD3	3:A:1804:PX4:C3	1.47	1.41
1:A:1207:ALA:CB	3:A:1804:PX4:H13	1.58	1.32
1:A:370:LYS:CD	3:A:1804:PX4:C3	2.28	1.11
1:A:370:LYS:HD3	3:A:1804:PX4:H5	1.31	1.08
1:A:968:VAL:O	1:A:995:GLN:O	1.72	1.07
1:A:966:THR:O	1:A:993:GLU:CB	2.03	1.07
1:A:1207:ALA:HB2	3:A:1804:PX4:H13	1.04	1.04
1:A:370:LYS:HD3	3:A:1804:PX4:H6	1.42	0.99
1:A:373:MET:HG2	3:A:1804:PX4:H11	1.42	0.98
1:A:1207:ALA:CB	3:A:1804:PX4:C5	2.43	0.95
1:A:1207:ALA:HB2	3:A:1804:PX4:C5	1.95	0.94
1:A:1386:LEU:HD11	1:A:1435:LEU:HD11	1.53	0.90
1:A:1207:ALA:HB3	3:A:1804:PX4:H13	1.54	0.90
1:A:373:MET:HG2	3:A:1804:PX4:C5	2.06	0.86
1:A:613:GLN:NE2	1:A:690:GLU:OE1	2.07	0.85
1:A:370:LYS:CD	3:A:1804:PX4:H7	2.15	0.76
1:A:1388:LYS:HE3	1:A:1390:TYR:HE1	1.50	0.76
1:A:1074:VAL:HG23	1:A:1075:SER:H	1.51	0.75
1:A:372:ASN:HB2	3:A:1804:PX4:H8	1.69	0.74
1:A:1207:ALA:HB3	3:A:1804:PX4:C5	2.14	0.74
1:A:1142:TRP:NE1	1:A:1325:GLU:OE1	2.20	0.74
1:A:558:GLU:HA	1:A:715:THR:HG22	1.70	0.73
1:A:1469:ASP:HB3	1:A:1513:ARG:HG3	1.69	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1012:ALA:HB2	1:A:1019:PHE:CE2	2.24	0.72
1:A:96:THR:HA	1:A:99:ARG:HG2	1.72	0.71
1:A:622:THR:HB	1:A:659:ASN:HA	1.70	0.71
1:A:711:LYS:O	1:A:713:ASP:N	2.20	0.71
1:A:376:ALA:HB1	1:A:1211:THR:HG21	1.71	0.71
1:A:977:GLN:HG3	1:A:978:LEU:H	1.55	0.71
1:A:1240:GLU:O	1:A:1240:GLU:HG2	1.90	0.70
1:A:1:MET:HG2	1:A:2:ALA:H	1.55	0.70
1:A:1292:GLN:HG3	1:A:1300:ALA:HB1	1.72	0.70
1:A:1372:SER:HB2	1:A:1375:SER:HB2	1.72	0.70
1:A:1149:ASP:OD1	1:A:1197:TYR:OH	2.08	0.70
1:A:370:LYS:CD	3:A:1804:PX4:H6	2.09	0.70
1:A:906:GLN:NE2	1:A:1488:GLU:OE2	2.24	0.69
1:A:399:MET:HE2	1:A:404:LYS:HG2	1.74	0.68
1:A:1239:GLU:HA	1:A:1239:GLU:OE1	1.92	0.67
1:A:121:ARG:HG3	1:A:122:TYR:CD2	2.29	0.67
1:A:1123:ARG:NH1	1:A:1203:PHE:O	2.27	0.67
1:A:1164:PHE:CD2	1:A:1172:PHE:HE2	2.13	0.67
1:A:20:ARG:NH2	1:A:620:ASN:O	2.28	0.67
1:A:838:ARG:NH1	1:A:855:PRO:O	2.27	0.67
1:A:194:ARG:HH12	1:A:204:PRO:HD3	1.59	0.66
1:A:372:ASN:CB	3:A:1804:PX4:H8	2.26	0.66
1:A:1439:ASP:OD2	1:A:1446:ARG:NH2	2.30	0.66
1:A:1161:LEU:HD13	1:A:1178:MET:HG3	1.80	0.64
1:A:1672:LYS:HG3	1:A:1673:TYR:CD2	2.33	0.64
1:A:1021:GLU:OE2	1:A:1021:GLU:HA	1.98	0.63
1:A:1012:ALA:HB2	1:A:1019:PHE:CD2	2.34	0.63
1:A:1388:LYS:HE3	1:A:1390:TYR:CE1	2.33	0.63
1:A:234:LEU:HD13	1:A:1065:LEU:HD23	1.81	0.62
1:A:1379:THR:OG1	1:A:1382:ILE:HD11	2.02	0.60
1:A:129:VAL:O	1:A:161:ARG:NH2	2.36	0.59
1:A:1388:LYS:HE2	1:A:1433:GLU:HG2	1.83	0.59
1:A:82:TYR:HE2	1:A:85:SER:HA	1.67	0.59
1:A:9:LEU:HD21	1:A:632:GLN:HB2	1.85	0.59
1:A:666:SER:HB3	1:A:668:GLY:H	1.68	0.58
1:A:711:LYS:HG3	1:A:712:SER:H	1.69	0.58
1:A:370:LYS:CE	3:A:1804:PX4:H6	2.33	0.57
1:A:206:TYR:O	1:A:212:LEU:HD23	2.05	0.56
1:A:1208:THR:O	1:A:1212:ARG:HG3	2.05	0.56
1:A:53:ASN:HA	1:A:249:PRO:HB3	1.86	0.56
1:A:1212:ARG:NH1	4:A:1807:POV:H25	2.20	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1511:LEU:O	1:A:1513:ARG:N	2.39	0.56
1:A:246:PRO:HB2	1:A:1009:ILE:HD11	1.87	0.56
1:A:332:LYS:H	1:A:335:VAL:CG2	2.19	0.55
1:A:376:ALA:HB2	4:A:1805:POV:C21	2.37	0.55
1:A:1357:GLU:HG3	1:A:1358:ASP:H	1.71	0.55
1:A:1226:LEU:O	1:A:1230:ILE:HG12	2.06	0.55
1:A:156:PHE:CE2	1:A:211:PHE:HD1	2.24	0.55
1:A:1588:VAL:HG12	1:A:1593:LYS:HG3	1.88	0.55
1:A:332:LYS:H	1:A:335:VAL:HG22	1.72	0.55
1:A:1114:THR:HG22	1:A:1210:TYR:HE2	1.72	0.55
1:A:1202:PHE:HD2	4:A:1807:POV:H29	1.71	0.55
1:A:241:THR:OG1	1:A:1039:ASN:OD1	2.25	0.54
1:A:1233:ILE:HG23	1:A:1234:PRO:HD2	1.88	0.54
1:A:368:PHE:CZ	1:A:377:PHE:HD2	2.26	0.54
1:A:368:PHE:HZ	1:A:377:PHE:HD2	1.54	0.54
1:A:1012:ALA:HB2	1:A:1019:PHE:HE2	1.70	0.54
1:A:771:GLU:O	1:A:775:GLN:HG3	2.08	0.54
1:A:981:GLN:O	1:A:985:HIS:ND1	2.42	0.53
1:A:1011:ARG:O	1:A:1015:GLU:HG2	2.09	0.53
1:A:1164:PHE:CD2	1:A:1172:PHE:CE2	2.97	0.53
1:A:1586:ILE:HG12	1:A:1594:CYS:SG	2.49	0.53
1:A:117:GLU:O	1:A:121:ARG:NE	2.40	0.53
1:A:729:LEU:HD12	1:A:729:LEU:O	2.09	0.53
1:A:1212:ARG:HH12	4:A:1807:POV:H25	1.72	0.53
1:A:762:LEU:HD13	1:A:822:ILE:HG12	1.90	0.52
1:A:925:MET:HG3	3:A:1804:PX4:H46	1.91	0.52
1:A:195:GLU:HG3	1:A:202:GLY:HA3	1.91	0.52
1:A:810:PHE:HD2	1:A:825:PHE:HE1	1.58	0.52
1:A:370:LYS:HD2	3:A:1804:PX4:H7	1.89	0.52
1:A:367:PHE:O	3:A:1803:PX4:H3	2.10	0.52
1:A:763:VAL:HB	1:A:823:ALA:O	2.10	0.52
1:A:1271:ARG:HB2	1:A:1291:TYR:CD2	2.45	0.52
1:A:843:VAL:HG12	1:A:843:VAL:O	2.11	0.51
1:A:1320:LEU:O	1:A:1324:ILE:HG12	2.11	0.51
1:A:157:SER:OG	1:A:158:TYR:N	2.43	0.51
1:A:801:GLU:N	1:A:801:GLU:OE2	2.43	0.51
1:A:1389:VAL:HG22	1:A:1397:LEU:HD12	1.92	0.51
1:A:933:PRO:HG2	1:A:1111:LEU:HD12	1.93	0.51
1:A:1171:ALA:HA	1:A:1268:TYR:CD2	2.46	0.50
1:A:332:LYS:O	1:A:335:VAL:HG22	2.11	0.50
1:A:1067:GLY:O	1:A:1069:HIS:N	2.44	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:977:GLN:HG3	1:A:978:LEU:N	2.26	0.50
1:A:148:LEU:HB3	1:A:1070:ALA:HB2	1.93	0.50
1:A:1074:VAL:HG23	1:A:1075:SER:N	2.24	0.50
1:A:1172:PHE:CD2	1:A:1181:THR:HG21	2.46	0.50
1:A:123:ASP:OD2	1:A:124:ASN:N	2.44	0.50
1:A:83:ILE:HG13	1:A:134:VAL:HG22	1.94	0.50
1:A:84:PRO:HG2	1:A:135:PHE:HB2	1.94	0.50
1:A:1066:CYS:SG	1:A:1067:GLY:N	2.85	0.49
1:A:1236:VAL:HG12	1:A:1237:LYS:HD3	1.94	0.49
1:A:1082:SER:OG	1:A:1085:GLN:HG3	2.12	0.49
1:A:123:ASP:OD2	1:A:125:CYS:N	2.41	0.49
1:A:219:ASP:O	1:A:223:MET:HG3	2.13	0.49
1:A:990:LEU:N	1:A:990:LEU:CD2	2.76	0.48
1:A:711:LYS:C	1:A:713:ASP:H	2.14	0.48
1:A:903:LEU:O	1:A:907:GLN:HG3	2.12	0.48
1:A:914:LYS:HG3	1:A:1479:MET:HE1	1.95	0.48
1:A:688:LEU:HD13	1:A:691:PRO:HB3	1.95	0.48
1:A:82:TYR:CE2	1:A:85:SER:HA	2.48	0.48
1:A:83:ILE:CG1	1:A:134:VAL:HG22	2.44	0.47
1:A:180:HIS:NE2	1:A:203:GLU:OE2	2.45	0.47
1:A:335:VAL:HG23	1:A:336:ALA:N	2.30	0.47
1:A:440:PRO:HB3	1:A:446:ASP:HA	1.96	0.47
1:A:982:LEU:HD21	1:A:1064:LEU:HD12	1.96	0.47
1:A:607:SER:O	1:A:683:SER:HB2	2.15	0.47
1:A:637:SER:HB2	1:A:640:LYS:HG2	1.96	0.47
1:A:990:LEU:N	1:A:990:LEU:HD23	2.29	0.47
1:A:1231:MET:HG3	1:A:1239:GLU:CB	2.45	0.47
1:A:753:LYS:HB2	1:A:753:LYS:HE2	1.67	0.47
1:A:151:LYS:HG2	1:A:1073:VAL:HG22	1.97	0.46
3:A:1804:PX4:H7	3:A:1804:PX4:H2	1.55	0.46
1:A:541:VAL:O	1:A:543:ASN:N	2.47	0.46
1:A:800:ARG:NH2	1:A:829:ILE:HD11	2.31	0.46
1:A:238:LEU:HG	1:A:1036:THR:O	2.15	0.46
1:A:325:LEU:O	1:A:329:VAL:HG13	2.15	0.46
1:A:684:LYS:O	1:A:714:ARG:HB2	2.15	0.46
1:A:810:PHE:CD2	1:A:825:PHE:HE1	2.33	0.46
1:A:1415:LEU:HD11	1:A:1584:LEU:HD22	1.97	0.46
1:A:1:MET:HG2	1:A:2:ALA:N	2.26	0.46
1:A:120:ILE:HG21	1:A:1080:PRO:HG2	1.98	0.46
1:A:1263:SER:OG	1:A:1301:PRO:O	2.33	0.46
1:A:1497:ASN:OD1	1:A:1498:THR:N	2.50	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1512:VAL:HG13	1:A:1512:VAL:O	2.15	0.45
3:A:1803:PX4:O8	3:A:1803:PX4:H14	2.17	0.45
1:A:218:VAL:O	1:A:222:ILE:HG13	2.17	0.45
1:A:431:GLY:O	1:A:433:GLN:HG2	2.16	0.45
1:A:1207:ALA:HB3	3:A:1804:PX4:H9	1.98	0.45
1:A:1006:GLU:CD	1:A:1006:GLU:H	2.19	0.45
1:A:686:LEU:HD21	1:A:688:LEU:HD23	1.99	0.45
3:A:1803:PX4:H69	3:A:1803:PX4:H62	1.49	0.45
1:A:977:GLN:CG	1:A:978:LEU:H	2.27	0.45
1:A:335:VAL:O	1:A:336:ALA:HB3	2.17	0.45
1:A:848:ASP:OD2	1:A:848:ASP:N	2.50	0.44
1:A:1011:ARG:O	1:A:1015:GLU:CG	2.65	0.44
1:A:1012:ALA:CB	1:A:1019:PHE:CD2	3.01	0.44
1:A:966:THR:C	1:A:993:GLU:CB	2.82	0.44
1:A:567:HIS:CE1	1:A:840:GLY:HA3	2.52	0.44
1:A:967:VAL:HA	1:A:993:GLU:CB	2.46	0.44
1:A:1120:VAL:HG22	1:A:1200:ASN:HB2	1.99	0.44
1:A:1207:ALA:HB3	3:A:1804:PX4:C4	2.48	0.44
1:A:152:TYR:CE2	1:A:1074:VAL:HG12	2.52	0.44
1:A:1642:ASP:OD1	1:A:1643:GLU:N	2.46	0.44
1:A:354:ALA:HA	1:A:357:THR:HG22	2.00	0.44
1:A:1231:MET:HG3	1:A:1239:GLU:CG	2.48	0.44
1:A:1233:ILE:CG2	1:A:1234:PRO:HD2	2.48	0.44
1:A:1638:SER:HB3	1:A:1650:TYR:CD1	2.52	0.44
1:A:1624:LEU:HD12	1:A:1624:LEU:HA	1.86	0.44
1:A:616:ILE:HG13	1:A:616:ILE:O	2.17	0.43
1:A:624:ALA:HB2	1:A:658:TRP:NE1	2.33	0.43
1:A:471:GLU:O	1:A:471:GLU:HG2	2.17	0.43
1:A:734:ALA:HB1	1:A:741:LEU:HD11	2.00	0.43
1:A:1117:ILE:HG23	1:A:1206:ALA:HB1	2.00	0.43
1:A:1154:LEU:O	1:A:1158:LEU:N	2.48	0.43
1:A:728:LEU:HD23	1:A:728:LEU:HA	1.83	0.43
1:A:922:GLU:OE2	3:A:1804:PX4:O8	2.37	0.43
1:A:1580:LEU:O	1:A:1580:LEU:HG	2.17	0.43
1:A:955:MET:HE2	1:A:1077:PHE:CD1	2.53	0.43
1:A:1219:LEU:O	1:A:1223:ALA:HB3	2.19	0.43
1:A:1651:HIS:CG	1:A:1651:HIS:O	2.70	0.43
1:A:1383:ILE:H	1:A:1383:ILE:HG13	1.65	0.43
1:A:695:MET:HE1	1:A:703:ILE:HD12	2.00	0.43
1:A:757:GLY:HA3	1:A:798:LEU:O	2.19	0.43
1:A:620:ASN:HA	1:A:661:ARG:HD2	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1233:ILE:HG13	1:A:1237:LYS:HE3	2.00	0.43
1:A:156:PHE:CE2	1:A:211:PHE:CD1	3.06	0.43
1:A:253:ASP:HB3	1:A:256:LEU:HD13	2.01	0.42
1:A:570:ALA:HB2	1:A:737:ALA:O	2.18	0.42
1:A:925:MET:CG	3:A:1804:PX4:H46	2.49	0.42
1:A:247:TYR:CD1	1:A:248:PRO:HD2	2.54	0.42
1:A:927:ALA:O	3:A:1803:PX4:H61	2.19	0.42
1:A:1472:THR:HG23	1:A:1511:LEU:H	1.85	0.42
1:A:1413:GLY:HA3	1:A:1584:LEU:HD23	2.02	0.42
1:A:958:LEU:HD13	1:A:1051:ALA:HB1	2.01	0.42
4:A:1807:POV:H29	4:A:1807:POV:H21B	1.41	0.42
1:A:535:LEU:HD12	1:A:535:LEU:HA	1.85	0.42
1:A:570:ALA:HB1	1:A:739:GLY:H	1.84	0.42
1:A:332:LYS:HB2	1:A:335:VAL:HG11	2.02	0.42
1:A:837:LEU:HD13	1:A:852:ILE:HG21	2.02	0.41
1:A:1245:LEU:HD23	1:A:1245:LEU:HA	1.85	0.41
1:A:1630:PHE:O	1:A:1634:THR:HG22	2.21	0.41
1:A:535:LEU:HA	1:A:588:GLY:HA3	2.03	0.41
1:A:155:ARG:O	1:A:1078:PRO:HA	2.20	0.41
1:A:634:LYS:NZ	1:A:677:ILE:O	2.50	0.41
1:A:1619:GLY:O	1:A:1622:GLU:N	2.46	0.41
1:A:212:LEU:HD12	1:A:1054:LEU:HD13	2.03	0.41
1:A:619:ASP:O	1:A:620:ASN:HB2	2.21	0.41
1:A:1614:LYS:HE2	1:A:1614:LYS:HB2	1.88	0.41
3:A:1803:PX4:H57	3:A:1803:PX4:H63	1.29	0.41
1:A:225:TYR:HD1	1:A:226:HIS:CE1	2.38	0.41
1:A:1300:ALA:HB3	1:A:1302:GLY:H	1.84	0.41
1:A:1318:LEU:HD23	1:A:1318:LEU:HA	1.86	0.41
1:A:1386:LEU:HD12	1:A:1386:LEU:HA	1.72	0.41
1:A:1569:ILE:HD12	1:A:1569:ILE:HG23	1.85	0.41
1:A:368:PHE:HZ	1:A:377:PHE:CD2	2.37	0.41
1:A:972:VAL:O	1:A:972:VAL:HG13	2.21	0.41
1:A:1084:LEU:HD23	1:A:1084:LEU:HA	1.94	0.41
1:A:1435:LEU:HD23	1:A:1448:SER:HB3	2.04	0.40
1:A:1576:GLU:O	1:A:1580:LEU:HD23	2.21	0.40
1:A:1630:PHE:HB2	1:A:1673:TYR:HE1	1.86	0.40
1:A:86:HIS:CD2	1:A:141:HIS:O	2.74	0.40
1:A:248:PRO:HG3	1:A:1012:ALA:O	2.21	0.40
1:A:977:GLN:O	1:A:978:LEU:HB2	2.21	0.40
1:A:752:GLN:HG3	1:A:1677:ASP:OD2	2.21	0.40
1:A:1012:ALA:CB	1:A:1019:PHE:HD2	2.34	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:603:GLN:OE1	1:A:603:GLN:HA	2.21	0.40
1:A:1042:PHE:CD1	1:A:1050:PRO:HG3	2.57	0.40
1:A:1154:LEU:HA	1:A:1157:SER:HB3	2.03	0.40
1:A:1368:ILE:HD11	1:A:1403:SER:OG	2.21	0.40
1:A:56:ILE:HG22	1:A:246:PRO:CB	2.51	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	1586/1748 (91%)	1363 (86%)	211 (13%)	12 (1%)	16	46

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	156	PHE
1	A	1234	PRO
1	A	712	SER
1	A	1075	SER
1	A	1238	LEU
1	A	995	GLN
1	A	1590	GLY
1	A	1021	GLU
1	A	432	ILE
1	A	1074	VAL
1	A	1512	VAL
1	A	337	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	1339/1486 (90%)	1328 (99%)	11 (1%)	79 87

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	19	LYS
1	A	114	LYS
1	A	489	MET
1	A	990	LEU
1	A	1019	PHE
1	A	1095	ARG
1	A	1232	ARG
1	A	1237	LYS
1	A	1238	LEU
1	A	1239	GLU
1	A	1240	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	1079	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](#)

7 ligands are modelled in this entry.

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	$\# Z > 2$
4	POV	A	1805	-	15,15,51	0.96	1 (6%)	14,14,59	0.86	0
4	POV	A	1807	-	15,15,51	0.96	1 (6%)	14,14,59	0.85	0
2	NAG	A	1801	1	14,14,15	0.30	0	17,19,21	0.37	0
4	POV	A	1806	-	15,15,51	0.97	1 (6%)	14,14,59	0.85	0
2	NAG	A	1802	1	14,14,15	0.50	0	17,19,21	0.46	0
3	PX4	A	1803	-	45,45,45	1.00	2 (4%)	51,53,53	1.01	3 (5%)
3	PX4	A	1804	-	45,45,45	1.01	2 (4%)	51,53,53	1.05	3 (5%)

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
4	POV	A	1805	-	-	8/13/13/55	-
4	POV	A	1807	-	-	6/13/13/55	-
2	NAG	A	1801	1	-	0/6/23/26	0/1/1/1
4	POV	A	1806	-	-	7/13/13/55	-
2	NAG	A	1802	1	-	2/6/23/26	0/1/1/1
3	PX4	A	1803	-	-	28/49/49/49	-
3	PX4	A	1804	-	-	24/49/49/49	-

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1804	PX4	O5-C9	4.29	1.45	1.33
3	A	1803	PX4	O5-C9	4.25	1.45	1.33
3	A	1804	PX4	O7-C23	4.08	1.45	1.34
3	A	1803	PX4	O7-C23	4.00	1.45	1.34
4	A	1806	POV	C29-C210	3.62	1.52	1.31
4	A	1805	POV	C29-C210	3.60	1.52	1.31
4	A	1807	POV	C29-C210	3.60	1.52	1.31

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1804	PX4	O7-C23-C24	4.05	120.23	111.50
3	A	1803	PX4	O7-C23-C24	3.93	119.97	111.50
3	A	1804	PX4	O5-C9-C10	2.63	120.16	111.91
3	A	1803	PX4	O5-C9-C10	2.59	120.02	111.91
3	A	1804	PX4	C7-O7-C23	-2.44	111.78	117.79
3	A	1803	PX4	C7-O7-C23	-2.22	112.33	117.79

There are no chirality outliers.

All (75) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	1803	PX4	C1-O3-P1-O1
3	A	1803	PX4	C1-O3-P1-O4
3	A	1803	PX4	C6-O4-P1-O1
3	A	1803	PX4	C6-O4-P1-O2
3	A	1803	PX4	C6-O4-P1-O3
3	A	1803	PX4	O3-C1-C2-N1
3	A	1804	PX4	C1-O3-P1-O4
3	A	1804	PX4	C6-O4-P1-O2
3	A	1804	PX4	C24-C23-O7-C7
3	A	1804	PX4	O8-C23-O7-C7
4	A	1805	POV	C211-C210-C29-C28
3	A	1803	PX4	C29-C30-C31-C32
3	A	1804	PX4	C16-C17-C18-C19
2	A	1802	NAG	O5-C5-C6-O6
3	A	1803	PX4	C23-C24-C25-C26
3	A	1804	PX4	C23-C24-C25-C26
3	A	1803	PX4	C32-C33-C34-C35
4	A	1806	POV	C211-C210-C29-C28
3	A	1804	PX4	C6-O4-P1-O3
3	A	1804	PX4	C10-C9-O5-C8

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Mol	Chain	Res	Type	Atoms
3	A	1804	PX4	C25-C26-C27-C28
3	A	1803	PX4	C24-C25-C26-C27
4	A	1807	POV	C23-C24-C25-C26
3	A	1803	PX4	C19-C20-C21-C22
3	A	1804	PX4	C28-C29-C30-C31
3	A	1803	PX4	C30-C31-C32-C33
3	A	1803	PX4	C28-C29-C30-C31
3	A	1804	PX4	O6-C9-O5-C8
3	A	1804	PX4	C27-C28-C29-C30
3	A	1803	PX4	C17-C18-C19-C20
3	A	1803	PX4	C12-C13-C14-C15
4	A	1806	POV	C212-C213-C214-C215
3	A	1804	PX4	C9-C10-C11-C12
3	A	1803	PX4	C26-C27-C28-C29
4	A	1807	POV	C22-C23-C24-C25
4	A	1807	POV	C210-C211-C212-C213
4	A	1805	POV	C24-C25-C26-C27
3	A	1804	PX4	C14-C15-C16-C17
4	A	1805	POV	C21-C22-C23-C24
4	A	1805	POV	C211-C212-C213-C214
3	A	1803	PX4	C27-C28-C29-C30
4	A	1807	POV	C27-C28-C29-C210
4	A	1805	POV	C210-C211-C212-C213
3	A	1804	PX4	C24-C25-C26-C27
3	A	1804	PX4	C13-C14-C15-C16
2	A	1802	NAG	C4-C5-C6-O6
4	A	1805	POV	C26-C27-C28-C29
3	A	1803	PX4	C33-C34-C35-C36
3	A	1804	PX4	C29-C30-C31-C32
3	A	1803	PX4	C10-C9-O5-C8
3	A	1803	PX4	C31-C32-C33-C34
4	A	1806	POV	C213-C214-C215-C216
3	A	1804	PX4	C15-C16-C17-C18
3	A	1803	PX4	C15-C16-C17-C18
3	A	1804	PX4	C26-C27-C28-C29
4	A	1806	POV	C21-C22-C23-C24
3	A	1803	PX4	O6-C9-O5-C8
3	A	1803	PX4	C1-O3-P1-O2
3	A	1804	PX4	C1-O3-P1-O1
3	A	1804	PX4	C6-O4-P1-O1
4	A	1805	POV	C22-C23-C24-C25
3	A	1803	PX4	C6-C7-O7-C23

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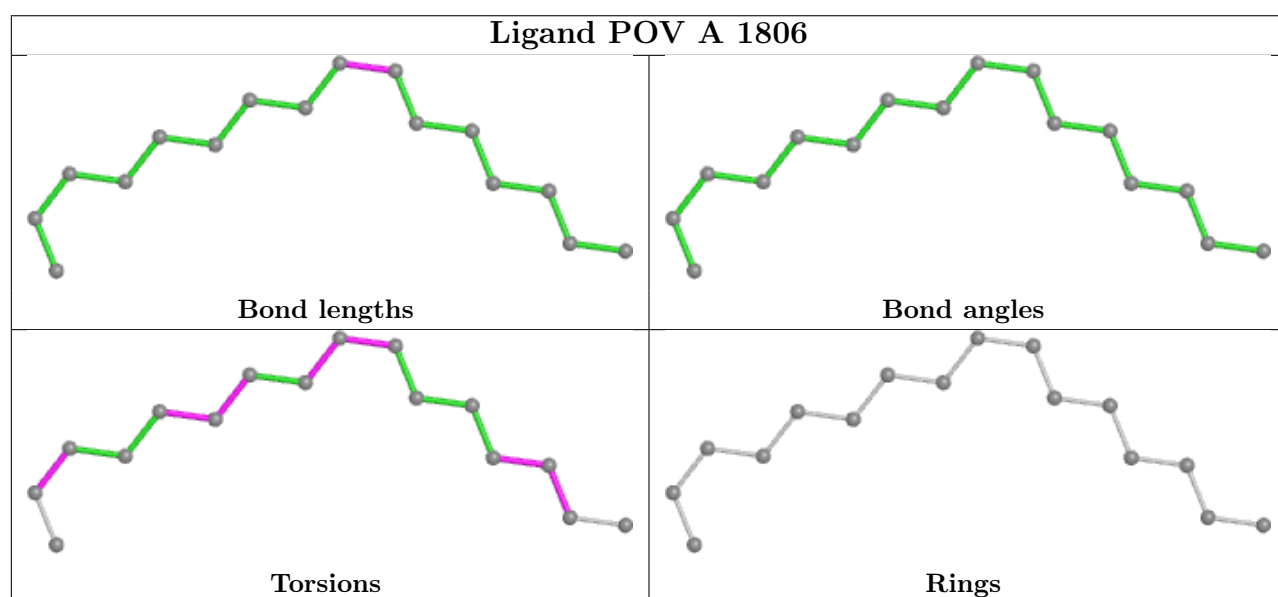
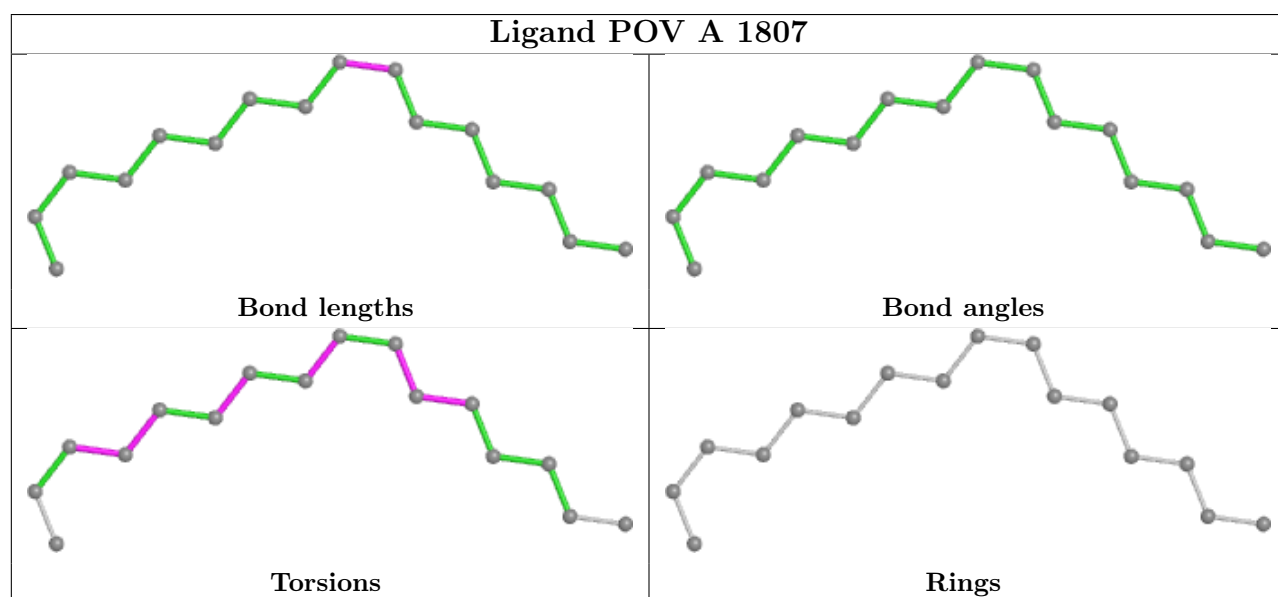
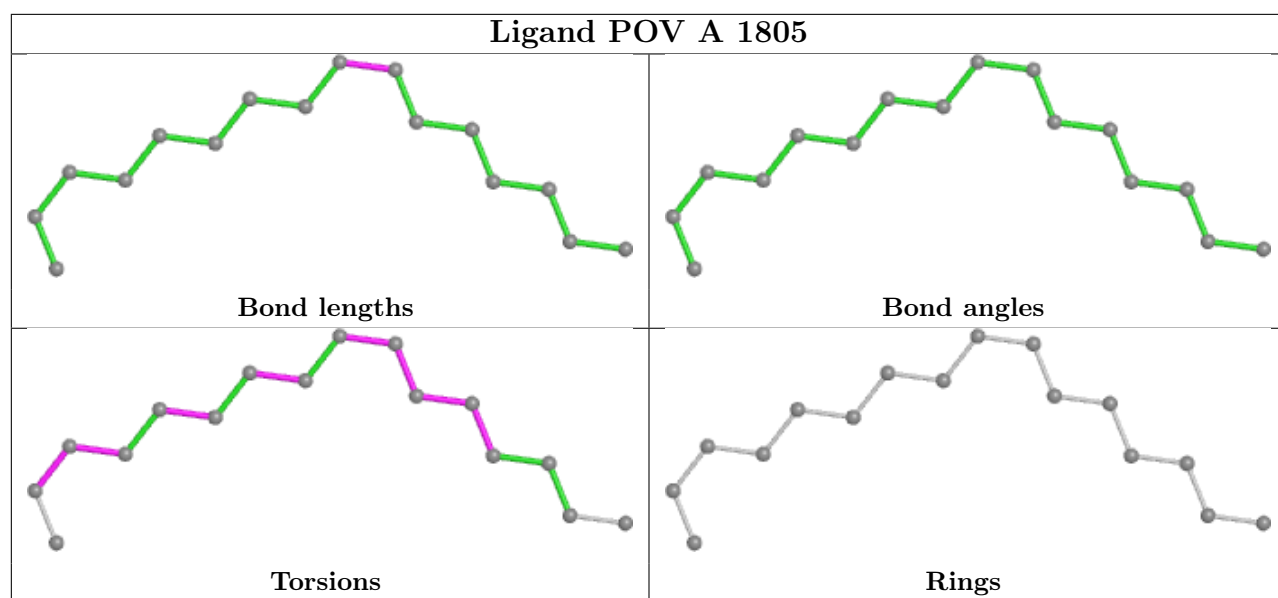
Mol	Chain	Res	Type	Atoms
4	A	1806	POV	C27-C28-C29-C210
4	A	1807	POV	C25-C26-C27-C28
3	A	1803	PX4	C10-C11-C12-C13
4	A	1806	POV	C24-C25-C26-C27
3	A	1803	PX4	C13-C14-C15-C16
3	A	1804	PX4	O7-C23-C24-C25
4	A	1806	POV	C25-C26-C27-C28
4	A	1805	POV	C29-C210-C211-C212
3	A	1804	PX4	C17-C18-C19-C20
3	A	1804	PX4	O8-C23-C24-C25
4	A	1807	POV	C29-C210-C211-C212
3	A	1803	PX4	O7-C23-C24-C25
3	A	1803	PX4	O8-C23-C24-C25

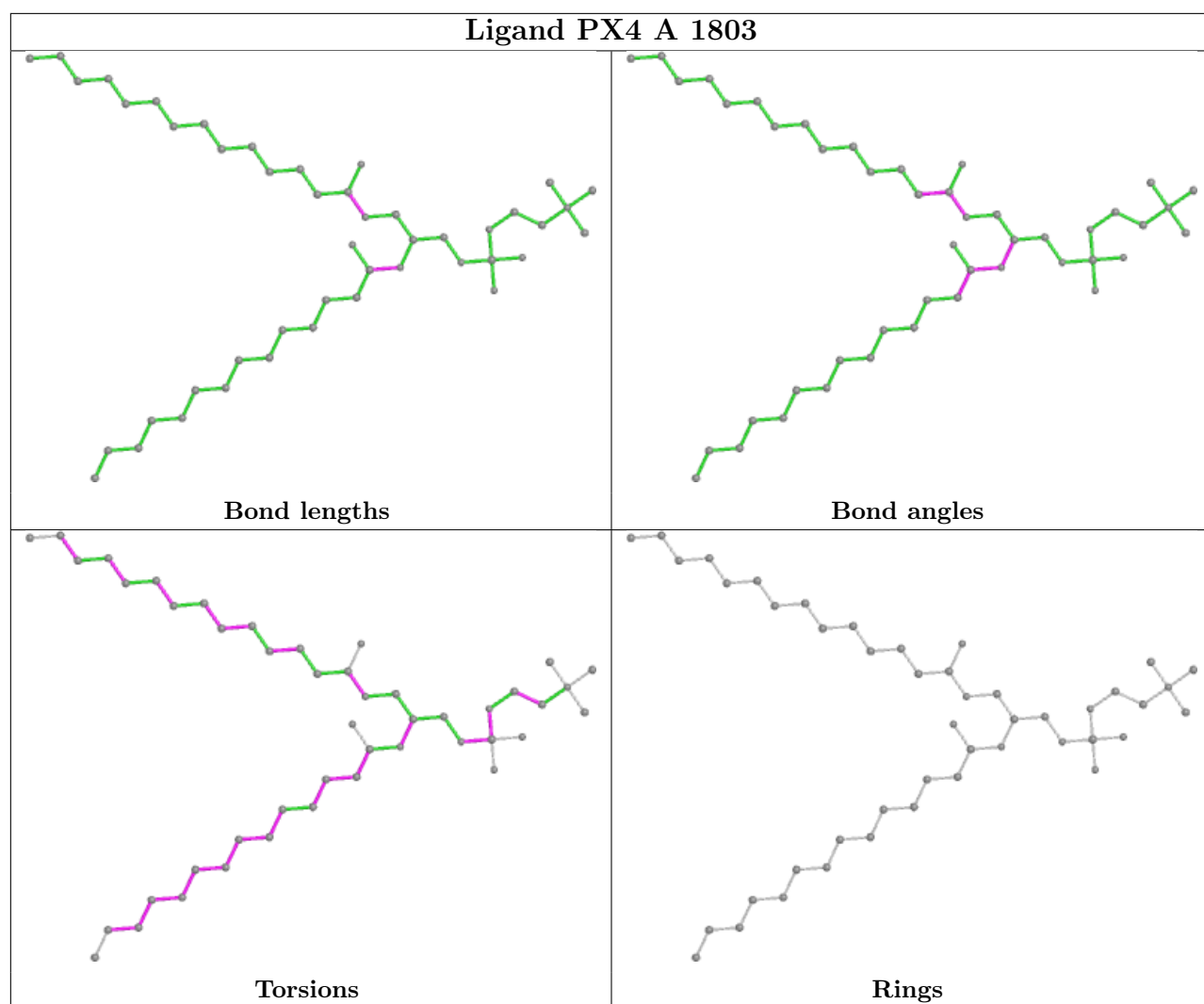
There are no ring outliers.

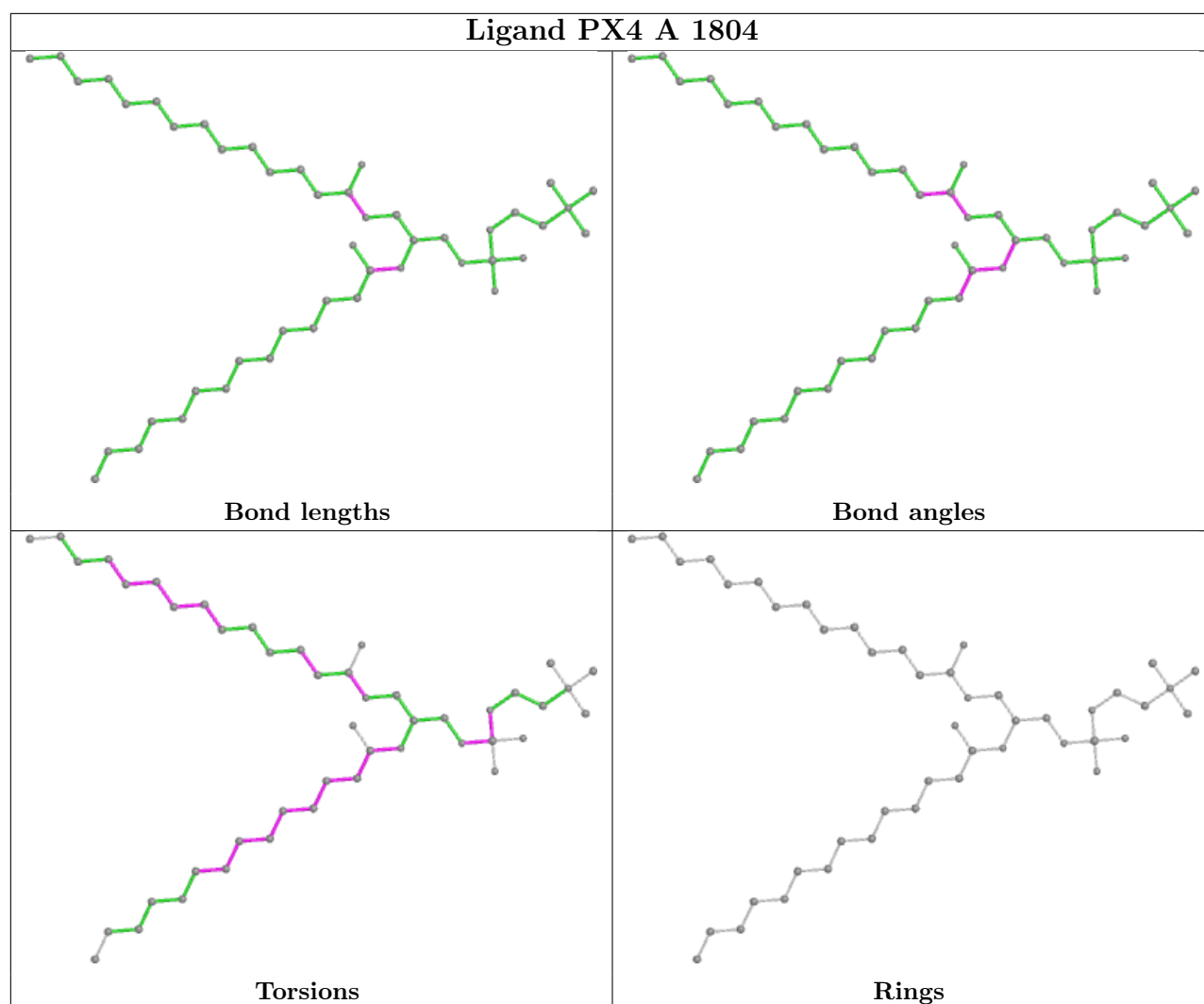
4 monomers are involved in 34 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	1805	POV	1	0
4	A	1807	POV	4	0
3	A	1803	PX4	5	0
3	A	1804	PX4	24	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.

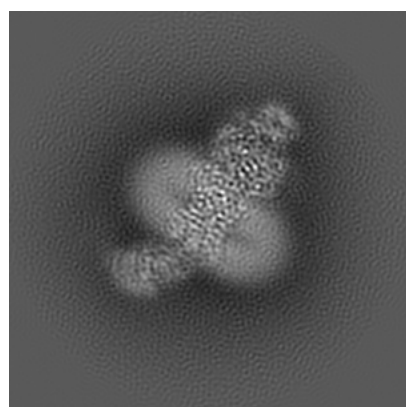
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-32233. These allow visual inspection of the internal detail of the map and identification of artifacts.

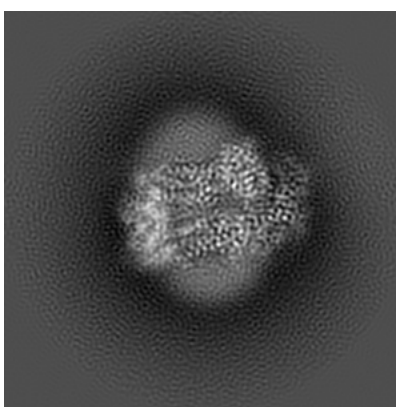
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

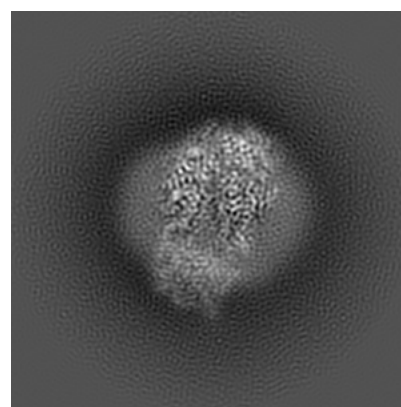
6.1.1 Primary map



X



Y

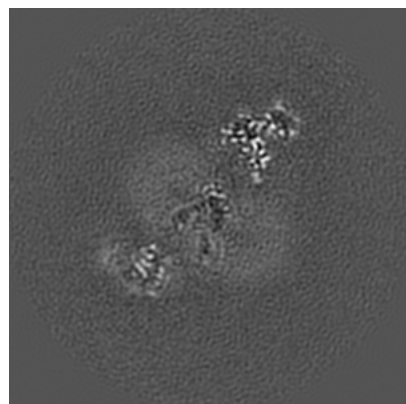


Z

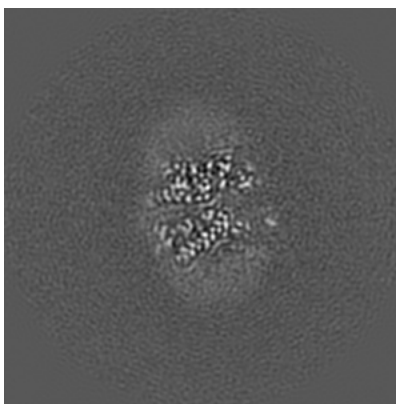
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

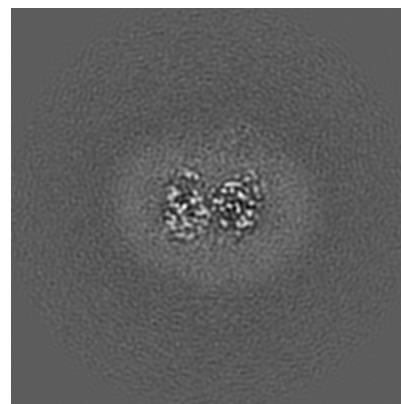
6.2.1 Primary map



X Index: 120



Y Index: 120

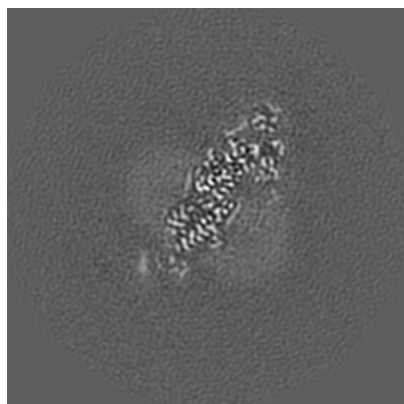


Z Index: 120

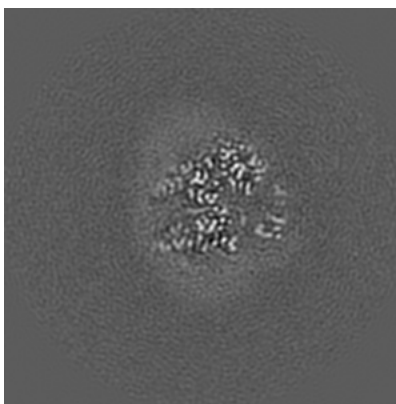
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

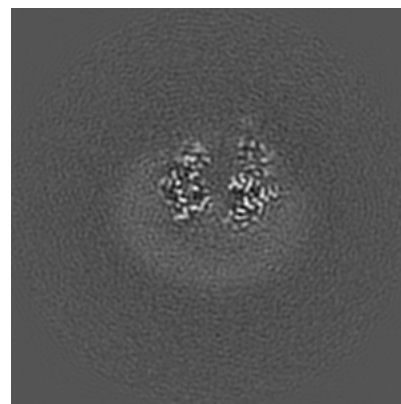
6.3.1 Primary map



X Index: 135



Y Index: 126

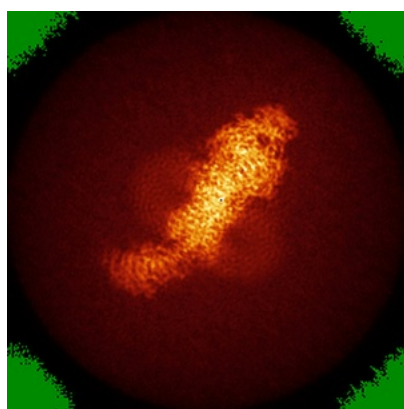


Z Index: 133

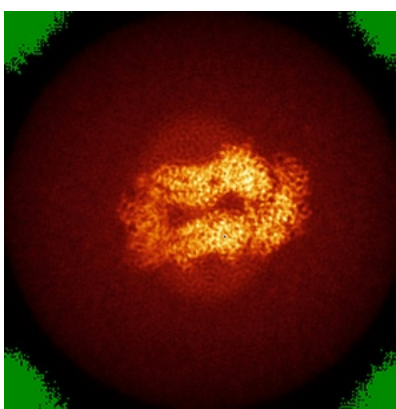
The images above show the largest variance slices of the map in three orthogonal directions.

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

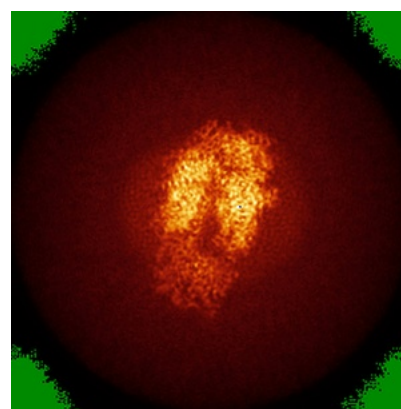
6.4.1 Primary map



X



Y



Z

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

6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.025. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

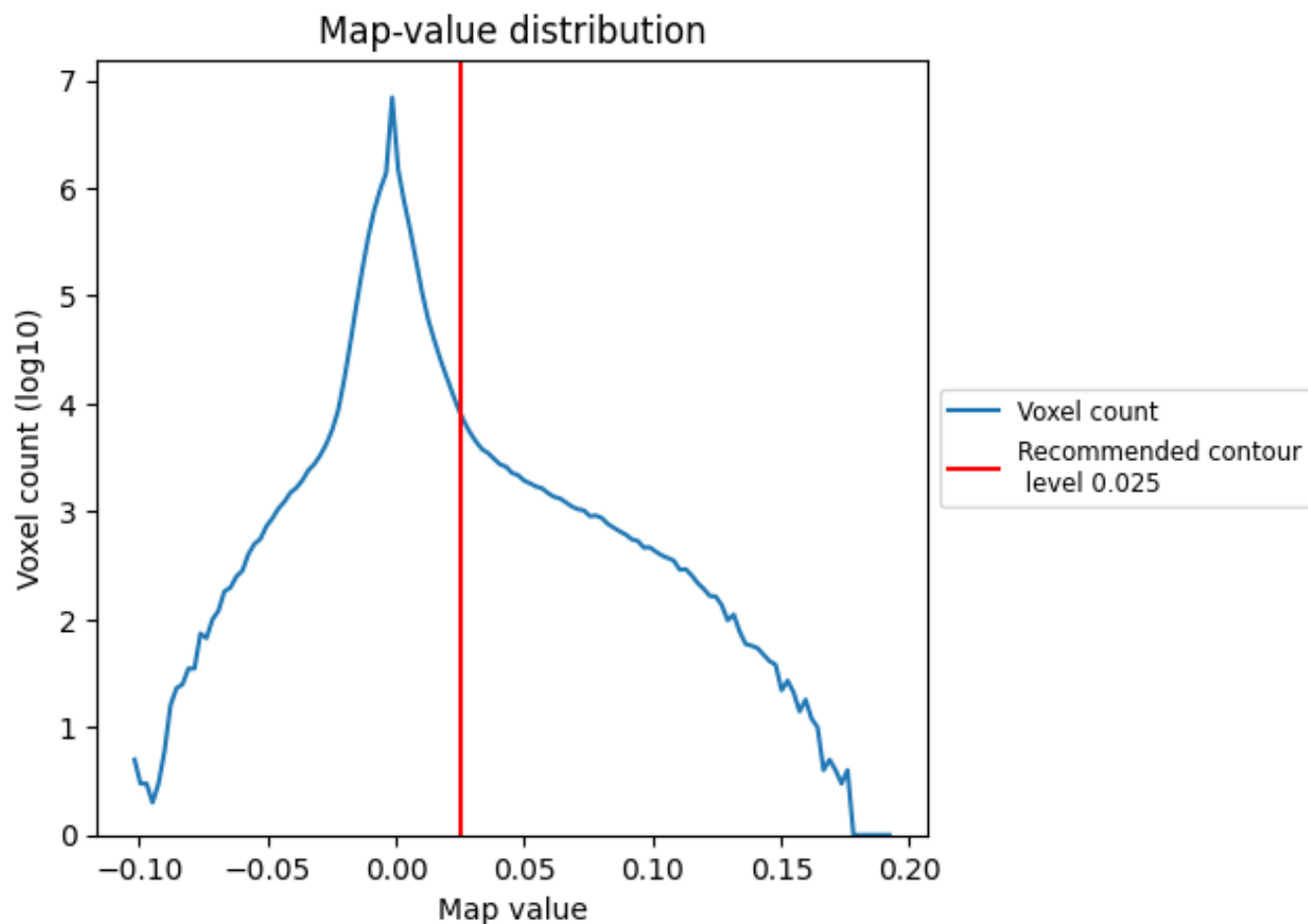
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

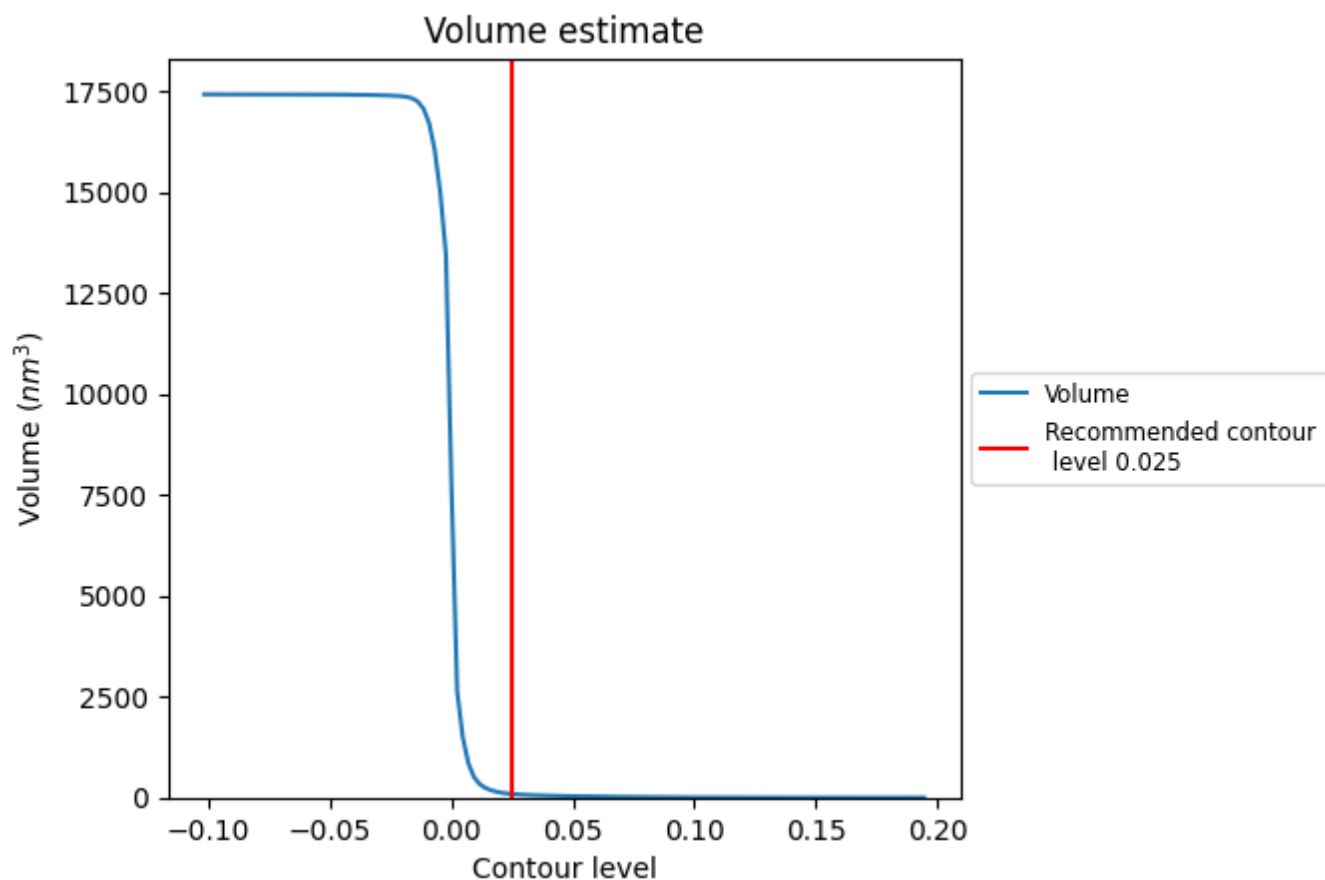
This section contains the results of statistical analysis of the map.

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.

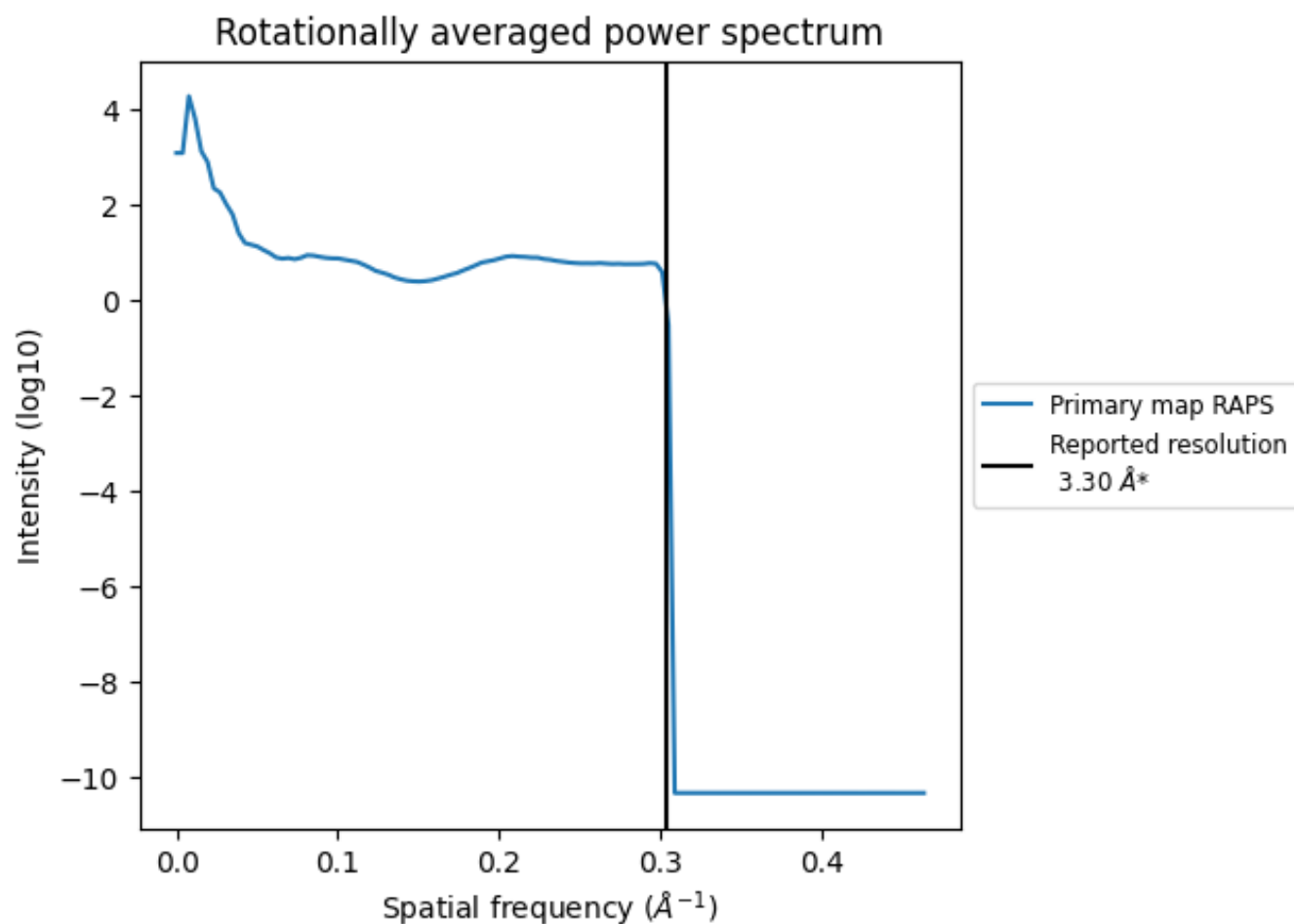
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 88 nm³; this corresponds to an approximate mass of 80 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 Å⁻¹

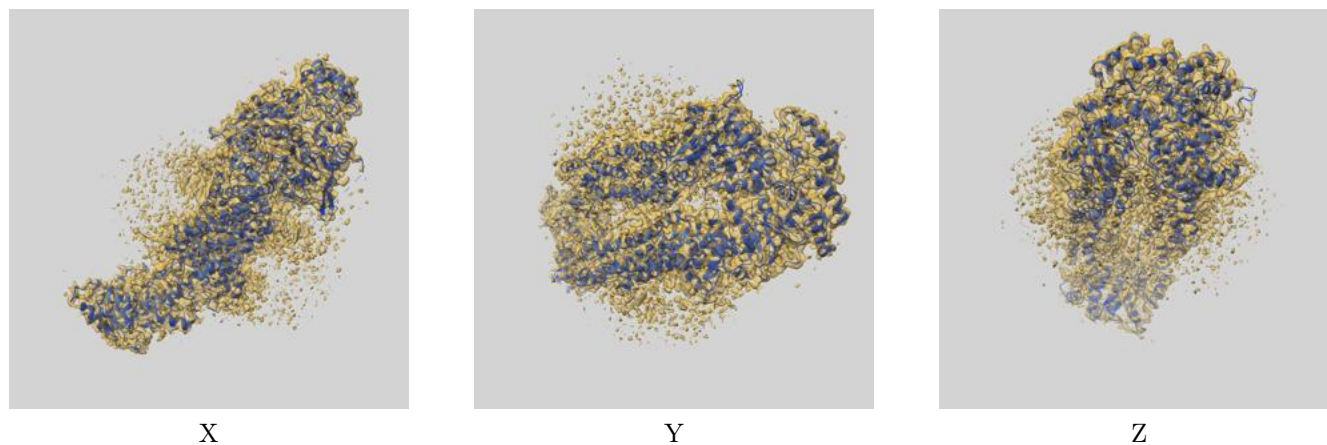
8 Fourier-Shell correlation

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

9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-32233 and PDB model 7W01. Per-residue inclusion information can be found in [section 3](#) on [page 7](#).

9.1 Map-model overlay [i](#)



The images above show the 3D surface view of the map at the recommended contour level 0.025 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](#)



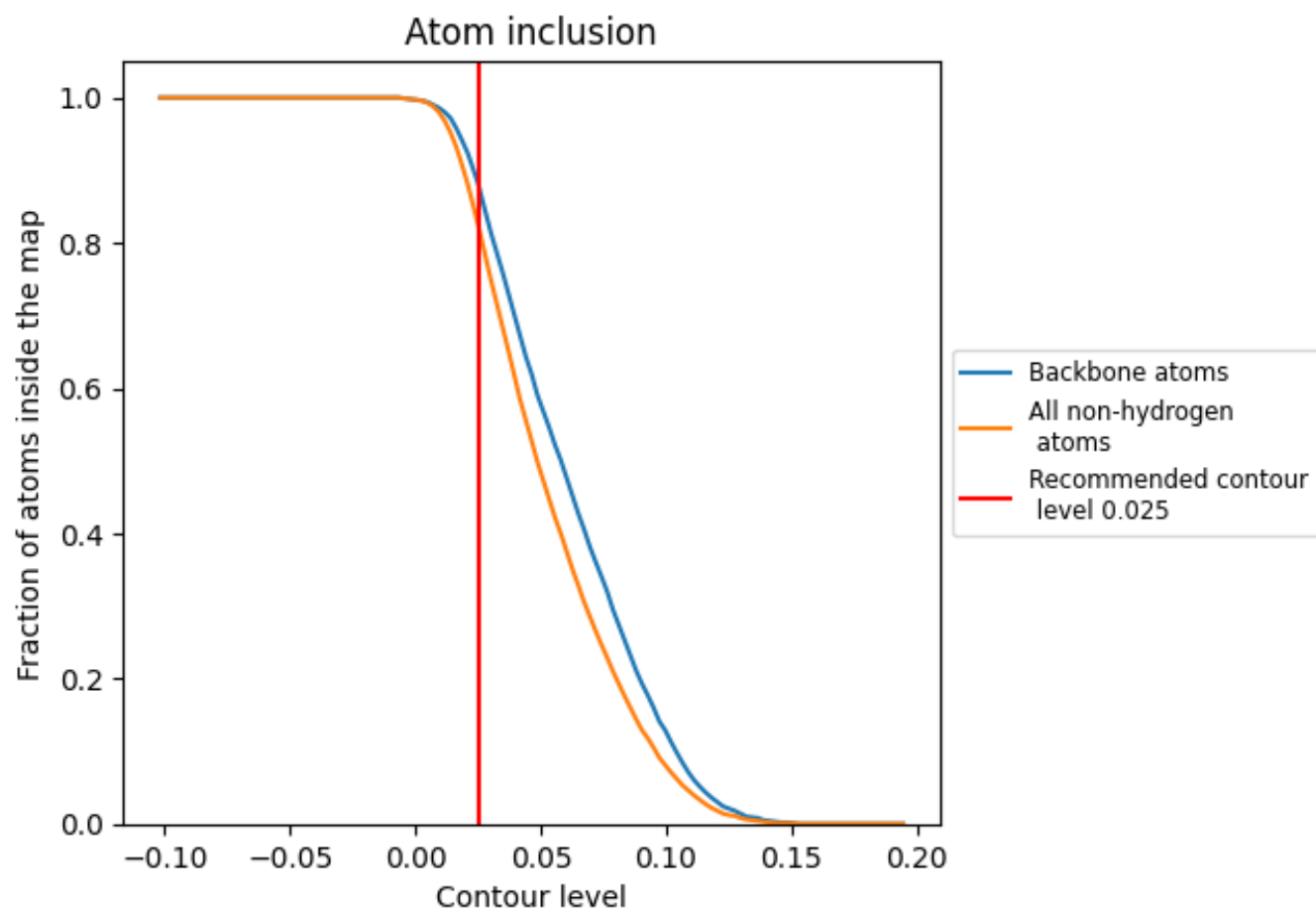
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.025).

9.4 Atom inclusion [i](#)



At the recommended contour level, 88% of all backbone atoms, 83% 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 (0.025) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	<div><div></div></div> 0.8260	<div><div></div></div> 0.5450
A	<div><div></div></div> 0.8260	<div><div></div></div> 0.5450

