Full wwPDB EM Validation Report ⓘ

May 27, 2025 – 10:17 AM JST

PDB ID	:	9IT0 / pdb_00009it0
EMDB ID	:	EMD-60853
Title	:	Liganded-state E.coli PatZ
Authors	:	Park, J.B.; Roh, S.H.
Deposited on	:	2024-07-19
Resolution	:	1.99 Å(reported)

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

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/EMValidationReportHelp
with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

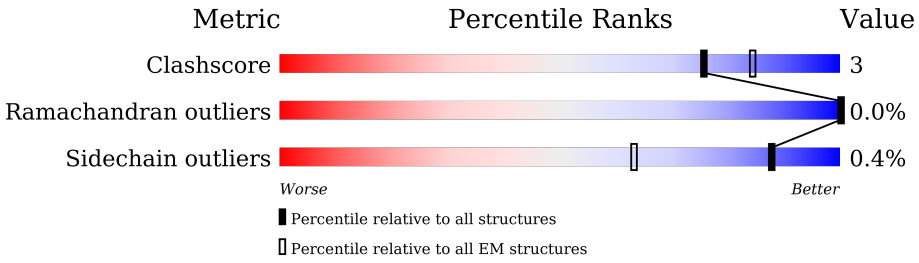
EMDB validation analysis	:	0.0.1.dev118
Mogul	:	1.8.5 (274361), CSD as541be (2020)
MolProbity	:	4-5-2 with Phenix2.0rc1
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
MapQ	:	1.9.13
Ideal geometry (proteins)	:	Engh & Huber (2001)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

1 Overall quality at a glance (i)

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

The reported resolution of this entry is 1.99 Å.

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



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

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

Mol	Chain	Length	Quality of chain
1	A	886	10% 93% 6%
1	B	886	10% 93% 6%
1	C	886	10% 94% 5%
1	D	886	10% 91% 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
2	ACO	A	902	X	-	-	-

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	ACO	B	901	X	-	-	-

2 Entry composition

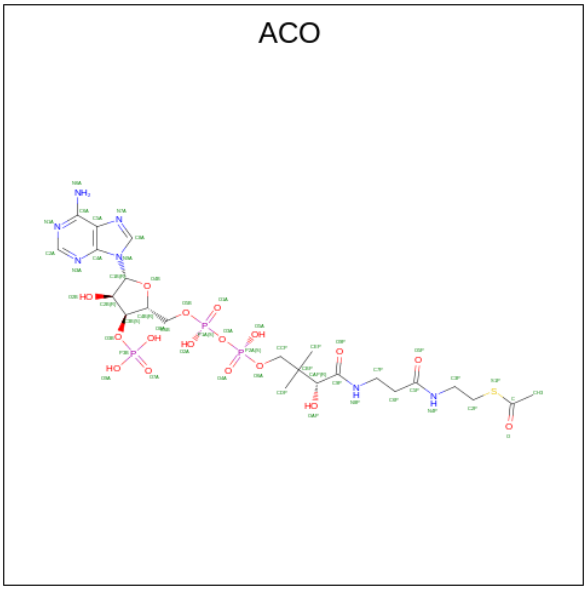
There are 4 unique types of molecules in this entry. The entry contains 56286 atoms, of which 27975 are hydrogens and 0 are deuteriums.

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 Protein acetyltransferase.

Mol	Chain	Residues	Atoms						AltConf	Trace
1	A	881	Total	C	H	N	O	S	0	0
			13770	4316	6926	1222	1275	31		
1	B	881	Total	C	H	N	O	S	0	0
			13770	4316	6926	1222	1275	31		
1	C	881	Total	C	H	N	O	S	0	0
			13770	4316	6926	1222	1275	31		
1	D	881	Total	C	H	N	O	S	0	0
			13770	4316	6926	1222	1275	31		

- Molecule 2 is ACETYL COENZYME *A (CCD ID: ACO) (formula: $C_{23}H_{38}N_7O_{17}P_3S$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms							AltConf
2	A	1	Total	C	H	N	O	P	S	0
			85	23	34	7	17	3	1	
2	A	1	Total	C	H	N	O	P	S	0
			84	23	33	7	17	3	1	

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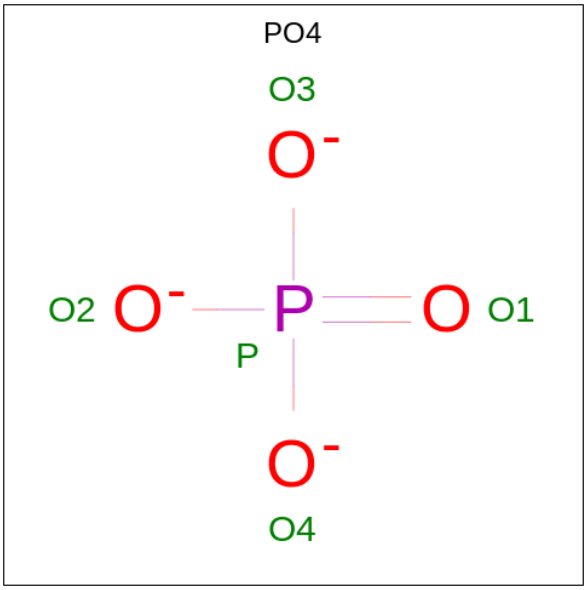
Mol	Chain	Residues	Atoms							AltConf
2	B	1	Total	C	H	N	O	P	S	0
			85	23	34	7	17	3	1	
2	B	1	Total	C	H	N	O	P	S	0
			85	23	34	7	17	3	1	
2	C	1	Total	C	H	N	O	P	S	0
			85	23	34	7	17	3	1	
2	C	1	Total	C	H	N	O	P	S	0
			85	23	34	7	17	3	1	
2	D	1	Total	C	H	N	O	P	S	0
			85	23	34	7	17	3	1	
2	D	1	Total	C	H	N	O	P	S	0
			85	23	34	7	17	3	1	

- Molecule 3 is PHOSPHATE ION (CCD ID: PO4) (formula: O_4P) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				AltConf
3	B	1	Total	O	P		0
			5	4	1		
3	B	1	Total	O	P		0
			5	4	1		
3	C	1	Total	O	P		0
			5	4	1		
3	C	1	Total	O	P		0
			5	4	1		

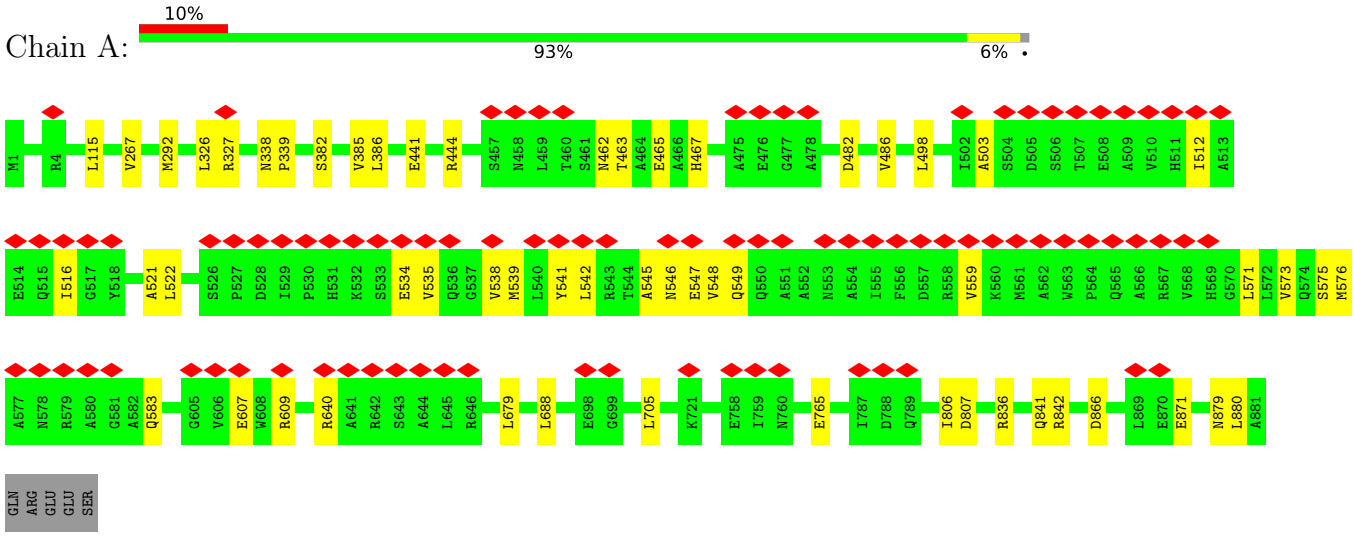
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	AltConf
4	A	106	Total O 106 106	0
4	B	110	Total O 110 110	0
4	C	121	Total O 121 121	0
4	D	170	Total O 170 170	0

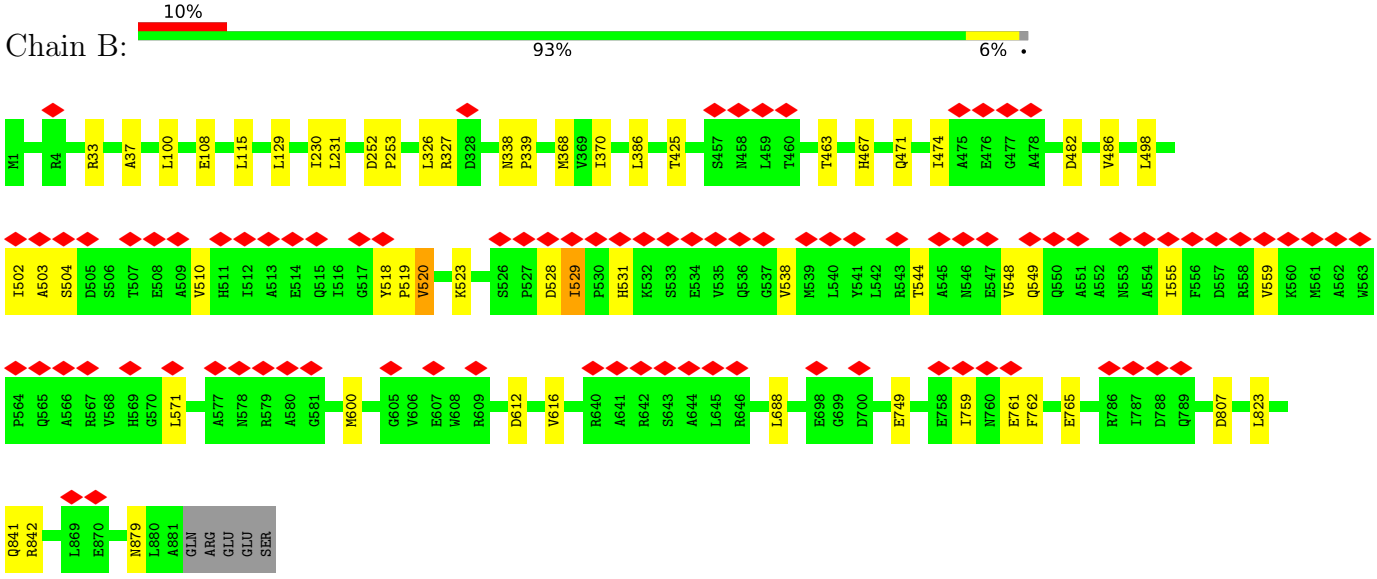
3 Residue-property plots (i)

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

• Molecule 1: Protein acetyltransferase

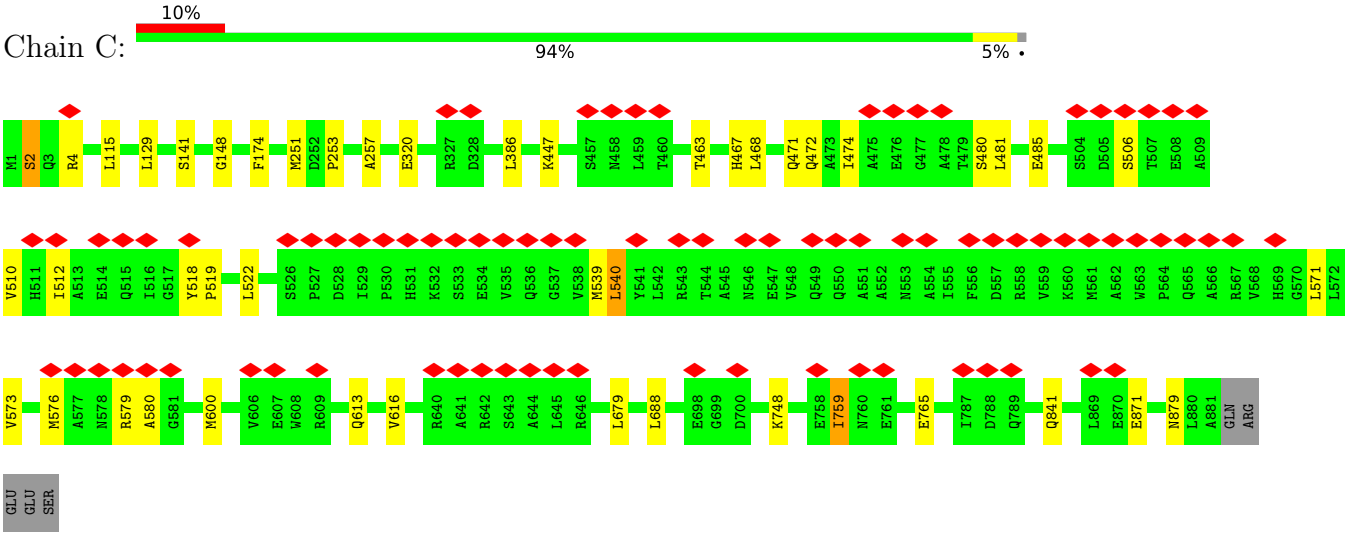


• Molecule 1: Protein acetyltransferase



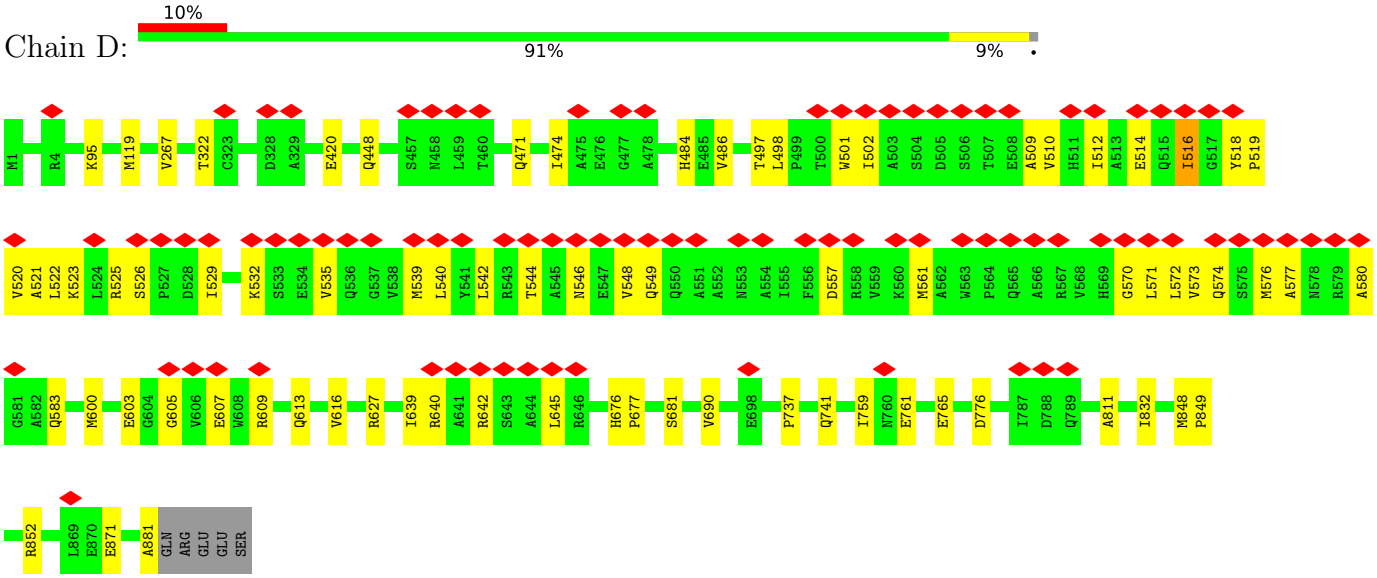
• Molecule 1: Protein acetyltransferase

Chain C:



• Molecule 1: Protein acetyltransferase

Chain D:



4 Experimental information (i)

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	5121944	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	NONE	Depositor
Microscope	TFS KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/Å^2$)	50	Depositor
Minimum defocus (nm)	1500	Depositor
Maximum defocus (nm)	8000	Depositor
Magnification	Not provided	
Image detector	FEI FALCON IV (4k x 4k)	Depositor
Maximum map value	1.315	Depositor
Minimum map value	0.000	Depositor
Average map value	0.010	Depositor
Map value standard deviation	0.020	Depositor
Recommended contour level	0.12	Depositor
Map size (Å)	340.47998, 340.47998, 340.47998	wwPDB
Map dimensions	448, 448, 448	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.75999993, 0.75999993, 0.75999993	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ACO, PO4

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	0.14	0/6976	0.31	0/9478
1	B	0.14	0/6976	0.33	0/9478
1	C	0.14	0/6976	0.32	0/9478
1	D	0.16	0/6976	0.34	0/9478
All	All	0.15	0/27904	0.32	0/37912

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	6844	6926	6926	38	0
1	B	6844	6926	6926	38	0
1	C	6844	6926	6926	36	0
1	D	6844	6926	6926	52	0
2	A	102	67	67	0	0
2	B	102	68	68	1	0
2	C	102	68	68	0	0
2	D	102	68	68	1	0
3	B	10	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	C	10	0	0	1	0
4	A	106	0	0	0	0
4	B	110	0	0	0	0
4	C	121	0	0	2	0
4	D	170	0	0	0	0
All	All	28311	27975	27975	163	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 3.

All (163) 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)
3:C:903:PO4:O2	4:C:1001:HOH:O	1.98	0.81
1:C:2:SER:OG	1:C:4:ARG:O	2.04	0.75
1:A:607:GLU:OE2	1:A:609:ARG:NH2	2.19	0.74
1:C:765:GLU:N	1:C:765:GLU:OE1	2.21	0.73
1:D:525:ARG:NH1	1:D:570:GLY:O	2.21	0.73
1:D:765:GLU:OE1	1:D:765:GLU:N	2.22	0.71
1:B:108:GLU:N	1:B:108:GLU:OE2	2.21	0.70
1:B:463:THR:HG22	1:B:467:HIS:CD2	2.26	0.70
1:D:849:PRO:O	1:D:852:ARG:NH2	2.23	0.70
1:B:765:GLU:N	1:B:765:GLU:OE1	2.24	0.70
1:C:871:GLU:OE1	1:C:871:GLU:O	2.10	0.69
1:A:765:GLU:N	1:A:765:GLU:OE1	2.25	0.69
1:C:141:SER:O	4:C:1002:HOH:O	2.10	0.69
1:D:516:ILE:HD11	1:D:520:VAL:HG13	1.75	0.68
1:A:539:MET:SD	1:A:576:MET:HE1	2.33	0.68
1:D:512:ILE:HG21	1:D:522:LEU:HD11	1.74	0.68
1:A:871:GLU:OE2	1:A:871:GLU:O	2.12	0.67
1:D:871:GLU:O	1:D:871:GLU:OE1	2.13	0.67
1:A:498:LEU:HD22	1:A:688:LEU:HD22	1.77	0.67
1:D:557:ASP:O	1:D:561:MET:HG2	1.94	0.67
1:B:529:ILE:HD12	1:B:531:HIS:O	1.94	0.66
1:D:676:HIS:HB3	1:D:677:PRO:HD3	1.78	0.65
1:A:546:ASN:OD1	1:A:547:GLU:N	2.30	0.65
1:D:642:ARG:NH1	1:D:645:LEU:O	2.29	0.65
1:C:539:MET:N	1:C:539:MET:HE3	2.11	0.65
1:A:880:LEU:HD13	1:A:880:LEU:O	1.96	0.64
1:D:523:LYS:HD3	1:D:572:LEU:HD11	1.79	0.64
1:D:502:ILE:HD12	1:D:572:LEU:HB3	1.80	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:600:MET:HG2	1:B:616:VAL:HG12	1.81	0.62
1:D:600:MET:HG2	1:D:616:VAL:HG12	1.81	0.62
1:D:519:PRO:HB2	1:D:576:MET:SD	2.40	0.62
1:A:462:ASN:ND2	1:A:465:GLU:OE2	2.33	0.61
1:D:526:SER:O	1:D:529:ILE:HG22	2.00	0.61
1:D:759:ILE:HD11	1:D:761:GLU:O	2.00	0.60
1:B:518:TYR:O	1:B:544:THR:OG1	2.20	0.60
1:A:534:GLU:OE2	1:A:559:VAL:HG12	2.01	0.60
1:A:535:VAL:HG12	1:A:538:VAL:HG23	1.85	0.59
1:D:832:ILE:HG23	1:D:881:ALA:O	2.02	0.58
1:D:516:ILE:HD11	1:D:520:VAL:CG1	2.33	0.58
1:D:681:SER:O	1:D:681:SER:OG	2.23	0.57
1:C:576:MET:HG3	1:C:576:MET:O	2.04	0.57
1:B:759:ILE:HD11	1:B:761:GLU:O	2.05	0.57
1:B:129:LEU:HD12	1:B:129:LEU:C	2.30	0.56
1:A:441:GLU:OE1	1:A:444:ARG:NH1	2.38	0.56
1:C:506:SER:O	1:C:510:VAL:HG23	2.05	0.55
1:D:521:ALA:HB3	1:D:574:GLN:HB2	1.89	0.55
1:A:583:GLN:NE2	1:A:640:ARG:O	2.39	0.55
1:B:338:ASN:HB2	1:B:339:PRO:HD3	1.89	0.54
1:A:521:ALA:HB1	1:A:538:VAL:O	2.08	0.54
1:A:866:ASP:O	1:A:866:ASP:OD1	2.26	0.53
1:C:251:MET:HG2	1:C:253:PRO:HD2	1.90	0.53
1:C:579:ARG:O	1:C:579:ARG:HG2	2.09	0.53
1:D:539:MET:HA	1:D:542:LEU:HD13	1.90	0.53
1:B:115:LEU:O	1:B:115:LEU:HD13	2.08	0.53
1:C:600:MET:HG2	1:C:616:VAL:HG12	1.91	0.53
1:B:100:LEU:HD11	2:B:901:ACO:H31	1.91	0.52
1:B:326:LEU:HD23	1:B:327:ARG:HE	1.74	0.52
1:C:129:LEU:C	1:C:129:LEU:HD12	2.34	0.52
1:A:338:ASN:HB3	1:A:339:PRO:HD3	1.91	0.52
1:B:612:ASP:N	1:B:612:ASP:OD1	2.42	0.52
1:B:504:SER:N	1:B:571:LEU:O	2.42	0.52
1:D:484:HIS:CD2	1:D:572:LEU:HD22	2.45	0.52
1:D:522:LEU:HD13	1:D:573:VAL:HA	1.92	0.51
1:B:498:LEU:HD22	1:B:688:LEU:HD22	1.92	0.51
1:A:545:ALA:O	1:A:549:GLN:NE2	2.43	0.51
1:C:512:ILE:HG21	1:C:573:VAL:HG11	1.93	0.51
1:A:386:LEU:C	1:A:386:LEU:HD23	2.36	0.51
1:A:503:ALA:O	1:A:571:LEU:N	2.43	0.51
1:C:447:LYS:O	1:C:447:LYS:HD3	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:605:GLY:O	1:D:607:GLU:N	2.44	0.50
1:A:836:ARG:HE	1:A:880:LEU:HD12	1.75	0.50
1:D:448:GLN:OE1	1:D:627:ARG:NH2	2.42	0.49
1:C:468:LEU:O	1:C:472:GLN:HG3	2.12	0.49
1:B:230:ILE:O	1:B:231:LEU:HD23	2.12	0.49
1:C:386:LEU:C	1:C:386:LEU:HD23	2.38	0.49
1:D:510:VAL:HG11	1:D:549:GLN:NE2	2.28	0.49
1:C:447:LYS:HD3	1:C:447:LYS:C	2.38	0.48
1:C:579:ARG:HA	1:D:540:LEU:HD11	1.95	0.48
1:D:516:ILE:HD12	1:D:518:TYR:H	1.79	0.48
1:B:463:THR:HG22	1:B:467:HIS:NE2	2.29	0.48
1:C:115:LEU:C	1:C:115:LEU:HD13	2.38	0.48
1:C:463:THR:HG22	1:C:467:HIS:CE1	2.48	0.48
1:A:541:TYR:CZ	1:A:576:MET:HE2	2.49	0.48
1:B:555:ILE:O	1:B:559:VAL:HG22	2.14	0.48
1:B:749:GLU:H	1:B:749:GLU:CD	2.22	0.48
1:D:532:LYS:O	1:D:535:VAL:HG22	2.14	0.48
1:A:542:LEU:H	1:A:542:LEU:HD22	1.80	0.47
1:B:115:LEU:HD13	1:B:115:LEU:C	2.39	0.47
1:A:679:LEU:HD12	1:A:688:LEU:HD21	1.96	0.47
1:A:115:LEU:C	1:A:115:LEU:HD13	2.39	0.47
1:D:609:ARG:N	1:D:613:GLN:OE1	2.46	0.47
1:A:516:ILE:HD12	1:A:575:SER:HB3	1.98	0.46
1:B:386:LEU:C	1:B:386:LEU:HD23	2.40	0.46
1:C:522:LEU:HD12	1:C:571:LEU:HD12	1.98	0.46
1:C:320:GLU:H	1:C:320:GLU:CD	2.23	0.46
1:A:338:ASN:OD1	1:A:338:ASN:C	2.59	0.46
1:D:486:VAL:HG21	1:D:690:VAL:HG13	1.98	0.46
1:D:516:ILE:HD12	1:D:518:TYR:N	2.31	0.46
1:D:603:GLU:H	1:D:603:GLU:CD	2.24	0.46
1:A:463:THR:HG22	1:A:467:HIS:CE1	2.51	0.45
1:C:679:LEU:HD22	1:C:688:LEU:CD1	2.47	0.45
1:B:368:MET:HE2	1:B:370:ILE:HD11	1.99	0.45
1:C:841:GLN:O	1:C:879:ASN:O	2.35	0.45
1:D:583:GLN:NE2	1:D:640:ARG:O	2.45	0.45
1:B:471:GLN:HA	1:B:474:ILE:HG22	1.99	0.45
1:C:580:ALA:HB2	1:D:580:ALA:HB1	1.97	0.45
1:A:292:MET:HE3	1:A:338:ASN:O	2.17	0.44
1:A:512:ILE:O	1:A:516:ILE:HG12	2.17	0.44
1:A:705:LEU:HD22	1:A:705:LEU:N	2.32	0.44
1:D:544:THR:OG1	1:D:546:ASN:OD1	2.30	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:555:ILE:O	1:B:559:VAL:HG13	2.17	0.44
1:A:841:GLN:O	1:A:879:ASN:O	2.36	0.44
1:B:33:ARG:HH12	1:B:37:ALA:HB2	1.83	0.44
1:B:510:VAL:HG12	1:B:548:VAL:CG2	2.47	0.44
1:B:523:LYS:HD3	1:B:538:VAL:HG11	1.99	0.44
1:B:807:ASP:OD1	1:B:842:ARG:NH1	2.51	0.44
1:D:811:ALA:HA	2:D:902:ACO:HH31	1.98	0.44
1:A:866:ASP:OD1	1:A:866:ASP:C	2.60	0.44
1:B:502:ILE:HG22	1:B:503:ALA:N	2.33	0.43
1:C:748:LYS:HE2	1:C:748:LYS:H	1.83	0.43
1:A:522:LEU:HD11	1:A:548:VAL:HG13	2.00	0.43
1:D:514:GLU:OE1	1:D:548:VAL:HG11	2.18	0.43
1:D:501:TRP:O	1:D:573:VAL:N	2.47	0.42
1:C:471:GLN:HA	1:C:474:ILE:HG22	2.01	0.42
1:C:679:LEU:HD22	1:C:688:LEU:HD11	2.00	0.42
1:B:252:ASP:N	1:B:253:PRO:HD2	2.34	0.42
1:C:148:GLY:N	1:C:174:PHE:O	2.53	0.42
1:C:759:ILE:HD12	1:C:759:ILE:N	2.35	0.42
1:B:544:THR:HG22	1:B:544:THR:O	2.19	0.42
1:D:529:ILE:O	1:D:529:ILE:HG23	2.19	0.42
1:B:841:GLN:O	1:B:879:ASN:O	2.38	0.42
1:B:482:ASP:O	1:B:486:VAL:HG22	2.20	0.42
1:D:523:LYS:HD3	1:D:572:LEU:CD1	2.49	0.42
1:D:497:THR:HG22	1:D:498:LEU:N	2.34	0.42
1:D:848:MET:HE3	1:D:848:MET:HB3	1.97	0.42
1:C:480:SER:O	1:C:481:LEU:HD23	2.20	0.42
1:D:471:GLN:HA	1:D:474:ILE:HG22	2.01	0.42
1:A:807:ASP:OD1	1:A:842:ARG:NH1	2.53	0.41
1:C:679:LEU:HD13	1:C:688:LEU:HD13	2.02	0.41
1:D:322:THR:HG22	1:D:322:THR:O	2.20	0.41
1:D:509:ALA:HB2	1:D:571:LEU:HD21	2.03	0.41
1:D:523:LYS:HE2	1:D:523:LYS:HB2	1.95	0.41
1:D:521:ALA:O	1:D:574:GLN:N	2.51	0.41
1:B:762:PHE:CD1	1:B:762:PHE:N	2.89	0.41
1:C:540:LEU:HB2	1:C:576:MET:SD	2.61	0.41
1:B:528:ASP:CG	1:B:529:ILE:H	2.28	0.41
1:A:482:ASP:O	1:A:486:VAL:HG22	2.20	0.41
1:B:519:PRO:O	1:B:520:VAL:HB	2.21	0.41
1:C:115:LEU:HD13	1:C:115:LEU:O	2.20	0.41
1:C:481:LEU:HD22	1:C:485:GLU:OE1	2.21	0.41
1:B:510:VAL:HG11	1:B:549:GLN:CD	2.46	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:257:ALA:HB2	1:C:616:VAL:HG23	2.03	0.40
1:A:326:LEU:HD23	1:A:327:ARG:HG3	2.02	0.40
1:D:510:VAL:HG11	1:D:549:GLN:CD	2.46	0.40
1:D:737:PRO:O	1:D:741:GLN:HG3	2.21	0.40
1:A:535:VAL:HG12	1:A:538:VAL:CG2	2.49	0.40
1:A:806:ILE:HG22	1:A:807:ASP:OD1	2.21	0.40
1:D:498:LEU:HD21	1:D:577:ALA:HB2	2.02	0.40
1:B:548:VAL:HG23	1:B:549:GLN:N	2.36	0.40
1:D:95:LYS:HA	1:D:119:MET:HG2	2.04	0.40
1:D:420:GLU:OE1	1:D:420:GLU:C	2.64	0.40
1:A:382:SER:O	1:A:385:VAL:HG22	2.22	0.40
1:C:518:TYR:N	1:C:519:PRO:HD2	2.36	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	879/886 (99%)	862 (98%)	17 (2%)	0	100	100
1	B	879/886 (99%)	854 (97%)	24 (3%)	1 (0%)	48	47
1	C	879/886 (99%)	857 (98%)	22 (2%)	0	100	100
1	D	879/886 (99%)	855 (97%)	24 (3%)	0	100	100
All	All	3516/3544 (99%)	3428 (98%)	87 (2%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	520	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	737/742 (99%)	735 (100%)	2 (0%)	91 94
1	B	737/742 (99%)	734 (100%)	3 (0%)	89 92
1	C	737/742 (99%)	733 (100%)	4 (0%)	86 90
1	D	737/742 (99%)	733 (100%)	4 (0%)	86 90
All	All	2948/2968 (99%)	2935 (100%)	13 (0%)	88 92

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	267	VAL
1	A	573	VAL
1	B	425	THR
1	B	529	ILE
1	B	823	LEU
1	C	2	SER
1	C	540	LEU
1	C	613	GLN
1	C	759	ILE
1	D	267	VAL
1	D	516	ILE
1	D	639	ILE
1	D	776	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (27) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	392	HIS
1	A	404	ASN
1	A	409	HIS
1	B	41	ASN
1	B	117	HIS
1	B	134	GLN

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Mol	Chain	Res	Type
1	B	282	HIS
1	B	362	GLN
1	B	565	GLN
1	B	868	GLN
1	C	41	ASN
1	C	117	HIS
1	C	268	GLN
1	C	392	HIS
1	C	409	HIS
1	C	412	GLN
1	C	515	GLN
1	C	569	HIS
1	C	574	GLN
1	C	655	GLN
1	D	268	GLN
1	D	282	HIS
1	D	392	HIS
1	D	409	HIS
1	D	591	HIS
1	D	655	GLN
1	D	712	HIS

5.3.3 RNA

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates

There are no oligosaccharides in this entry.

5.6 Ligand geometry

12 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	$\#\|Z\| > 2$
2	ACO	D	901	-	45,53,53	0.49	0	56,79,79	0.91	6 (10%)
3	PO4	B	904	-	4,4,4	0.93	0	6,6,6	0.43	0
3	PO4	C	903	-	4,4,4	0.93	0	6,6,6	0.46	0
3	PO4	C	904	-	4,4,4	0.93	0	6,6,6	0.49	0
2	ACO	B	901	-	45,53,53	0.49	0	56,79,79	1.11	6 (10%)
2	ACO	C	901	-	45,53,53	0.48	0	56,79,79	0.92	6 (10%)
2	ACO	A	902	-	45,53,53	0.50	0	56,79,79	1.11	6 (10%)
2	ACO	C	902	-	45,53,53	0.50	0	56,79,79	0.92	6 (10%)
2	ACO	B	902	-	45,53,53	0.49	0	56,79,79	0.91	6 (10%)
2	ACO	D	902	-	45,53,53	0.48	0	56,79,79	0.90	5 (8%)
2	ACO	A	901	-	45,53,53	0.50	0	56,79,79	0.92	6 (10%)
3	PO4	B	903	-	4,4,4	0.92	0	6,6,6	0.43	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
2	ACO	D	901	-	-	10/47/67/67	0/3/3/3
2	ACO	B	901	-	1/1/14/14	17/47/67/67	0/3/3/3
2	ACO	C	901	-	-	2/47/67/67	0/3/3/3
2	ACO	A	902	-	1/1/14/14	14/47/67/67	0/3/3/3
2	ACO	C	902	-	-	7/47/67/67	0/3/3/3
2	ACO	B	902	-	-	5/47/67/67	0/3/3/3
2	ACO	D	902	-	-	6/47/67/67	0/3/3/3
2	ACO	A	901	-	-	9/47/67/67	0/3/3/3

There are no bond length outliers.

All (47) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	902	ACO	OAP-CAP-CBP	5.21	122.53	110.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	901	ACO	OAP-CAP-CBP	5.07	122.18	110.25
2	C	901	ACO	O4B-C1B-C2B	-2.45	103.34	106.93
2	C	901	ACO	P2A-O6A-CCP	-2.44	107.51	121.56
2	B	901	ACO	O4B-C1B-C2B	-2.40	103.42	106.93
2	A	901	ACO	OAP-CAP-CBP	-2.39	104.64	110.25
2	D	901	ACO	OAP-CAP-CBP	-2.38	104.65	110.25
2	B	902	ACO	OAP-CAP-CBP	-2.36	104.69	110.25
2	C	902	ACO	OAP-CAP-CBP	-2.36	104.70	110.25
2	B	901	ACO	P2A-O6A-CCP	-2.34	108.09	121.56
2	C	901	ACO	OAP-CAP-CBP	-2.32	104.78	110.25
2	D	901	ACO	P2A-O6A-CCP	-2.32	108.19	121.56
2	B	901	ACO	O2B-C2B-C3B	-2.31	104.59	111.17
2	C	902	ACO	O2B-C2B-C3B	-2.31	104.60	111.17
2	D	901	ACO	O4B-C1B-C2B	-2.30	103.56	106.93
2	A	901	ACO	O4B-C1B-C2B	-2.30	103.56	106.93
2	C	901	ACO	O3B-C3B-C2B	-2.30	103.34	111.68
2	D	902	ACO	C5A-C6A-N6A	2.30	123.85	120.35
2	A	901	ACO	P2A-O6A-CCP	-2.29	108.33	121.56
2	C	902	ACO	C5A-C6A-N6A	2.29	123.83	120.35
2	B	901	ACO	C5A-C6A-N6A	2.29	123.83	120.35
2	B	902	ACO	C5A-C6A-N6A	2.29	123.83	120.35
2	A	901	ACO	C5A-C6A-N6A	2.28	123.82	120.35
2	C	901	ACO	C5A-C6A-N6A	2.28	123.81	120.35
2	D	901	ACO	C5A-C6A-N6A	2.28	123.81	120.35
2	A	902	ACO	C5A-C6A-N6A	2.28	123.81	120.35
2	B	901	ACO	O3B-C3B-C2B	-2.27	103.45	111.68
2	A	902	ACO	O4B-C1B-C2B	-2.26	103.63	106.93
2	D	902	ACO	O2B-C2B-C3B	-2.25	104.77	111.17
2	A	902	ACO	O3B-C3B-C2B	-2.25	103.53	111.68
2	B	902	ACO	O2B-C2B-C3B	-2.25	104.78	111.17
2	D	902	ACO	OAP-CAP-CBP	-2.24	104.97	110.25
2	C	901	ACO	O2B-C2B-C3B	-2.24	104.81	111.17
2	B	902	ACO	O4B-C1B-C2B	-2.24	103.66	106.93
2	B	902	ACO	O3B-C3B-C2B	-2.23	103.59	111.68
2	A	902	ACO	O2B-C2B-C3B	-2.23	104.84	111.17
2	D	902	ACO	O3B-C3B-C2B	-2.22	103.65	111.68
2	D	901	ACO	O3B-C3B-C2B	-2.20	103.70	111.68
2	A	901	ACO	O3B-C3B-C2B	-2.20	103.71	111.68
2	C	902	ACO	O3B-C3B-C2B	-2.19	103.74	111.68
2	A	901	ACO	O2B-C2B-C3B	-2.16	105.04	111.17
2	B	902	ACO	P2A-O6A-CCP	-2.15	109.18	121.56
2	D	902	ACO	O4B-C1B-C2B	-2.13	103.81	106.93

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	C	902	ACO	O4B-C1B-C2B	-2.12	103.83	106.93
2	D	901	ACO	O2B-C2B-C3B	-2.10	105.19	111.17
2	C	902	ACO	P2A-O6A-CCP	-2.06	109.70	121.56
2	A	902	ACO	P2A-O6A-CCP	-2.02	109.91	121.56

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	A	902	ACO	CAP
2	B	901	ACO	CAP

All (70) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	901	ACO	C5P-C6P-C7P-N8P
2	A	902	ACO	C5B-O5B-P1A-O1A
2	A	902	ACO	C5B-O5B-P1A-O2A
2	A	902	ACO	OAP-CAP-CBP-CCP
2	A	902	ACO	C9P-CAP-CBP-CCP
2	A	902	ACO	OAP-CAP-CBP-CDP
2	A	902	ACO	OAP-CAP-CBP-CEP
2	A	902	ACO	C9P-CAP-CBP-CEP
2	B	901	ACO	C3B-O3B-P3B-O9A
2	B	901	ACO	OAP-CAP-CBP-CCP
2	B	901	ACO	C9P-CAP-CBP-CCP
2	B	901	ACO	OAP-CAP-CBP-CDP
2	B	901	ACO	C9P-CAP-CBP-CDP
2	B	901	ACO	OAP-CAP-CBP-CEP
2	B	901	ACO	C9P-CAP-CBP-CEP
2	B	901	ACO	O-C-S1P-C2P
2	B	901	ACO	CH3-C-S1P-C2P
2	B	902	ACO	C5B-O5B-P1A-O1A
2	C	901	ACO	C3B-O3B-P3B-O8A
2	C	902	ACO	C5B-O5B-P1A-O1A
2	C	902	ACO	C5B-O5B-P1A-O2A
2	C	902	ACO	C3P-C2P-S1P-C
2	C	902	ACO	O-C-S1P-C2P
2	C	902	ACO	CH3-C-S1P-C2P
2	D	901	ACO	C5P-C6P-C7P-N8P
2	D	901	ACO	S1P-C2P-C3P-N4P
2	D	901	ACO	O-C-S1P-C2P
2	D	901	ACO	CH3-C-S1P-C2P

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Mol	Chain	Res	Type	Atoms
2	D	902	ACO	C5B-O5B-P1A-O1A
2	D	902	ACO	C5B-O5B-P1A-O2A
2	D	902	ACO	O-C-S1P-C2P
2	D	902	ACO	CH3-C-S1P-C2P
2	A	901	ACO	C6P-C5P-N4P-C3P
2	B	901	ACO	C6P-C5P-N4P-C3P
2	D	901	ACO	C6P-C5P-N4P-C3P
2	A	901	ACO	O5P-C5P-N4P-C3P
2	B	901	ACO	O5P-C5P-N4P-C3P
2	D	901	ACO	O5P-C5P-N4P-C3P
2	A	901	ACO	O-C-S1P-C2P
2	A	902	ACO	O9P-C9P-CAP-OAP
2	D	902	ACO	C6P-C7P-N8P-C9P
2	C	902	ACO	C6P-C7P-N8P-C9P
2	A	901	ACO	CH3-C-S1P-C2P
2	B	901	ACO	C3B-O3B-P3B-O7A
2	C	901	ACO	C3B-O3B-P3B-O7A
2	B	902	ACO	C5B-O5B-P1A-O3A
2	C	902	ACO	C5B-O5B-P1A-O3A
2	D	902	ACO	C5B-O5B-P1A-O3A
2	B	902	ACO	C5B-O5B-P1A-O2A
2	D	901	ACO	C2P-C3P-N4P-C5P
2	B	901	ACO	O9P-C9P-CAP-OAP
2	A	902	ACO	O4B-C4B-C5B-O5B
2	B	902	ACO	C6P-C7P-N8P-C9P
2	A	901	ACO	P2A-O3A-P1A-O2A
2	B	901	ACO	P2A-O3A-P1A-O2A
2	B	902	ACO	P1A-O3A-P2A-O5A
2	D	901	ACO	P2A-O3A-P1A-O2A
2	A	902	ACO	CEP-CBP-CCP-O6A
2	A	901	ACO	C3P-C2P-S1P-C
2	A	901	ACO	P2A-O3A-P1A-O1A
2	B	901	ACO	N8P-C9P-CAP-OAP
2	A	902	ACO	CDP-CBP-CCP-O6A
2	B	901	ACO	C2P-C3P-N4P-C5P
2	A	902	ACO	C9P-CAP-CBP-CDP
2	A	902	ACO	C6P-C7P-N8P-C9P
2	A	902	ACO	C5B-O5B-P1A-O3A
2	B	901	ACO	P2A-O3A-P1A-O1A
2	A	901	ACO	C5B-O5B-P1A-O1A
2	D	901	ACO	C5B-O5B-P1A-O1A
2	D	901	ACO	O5P-C5P-C6P-C7P

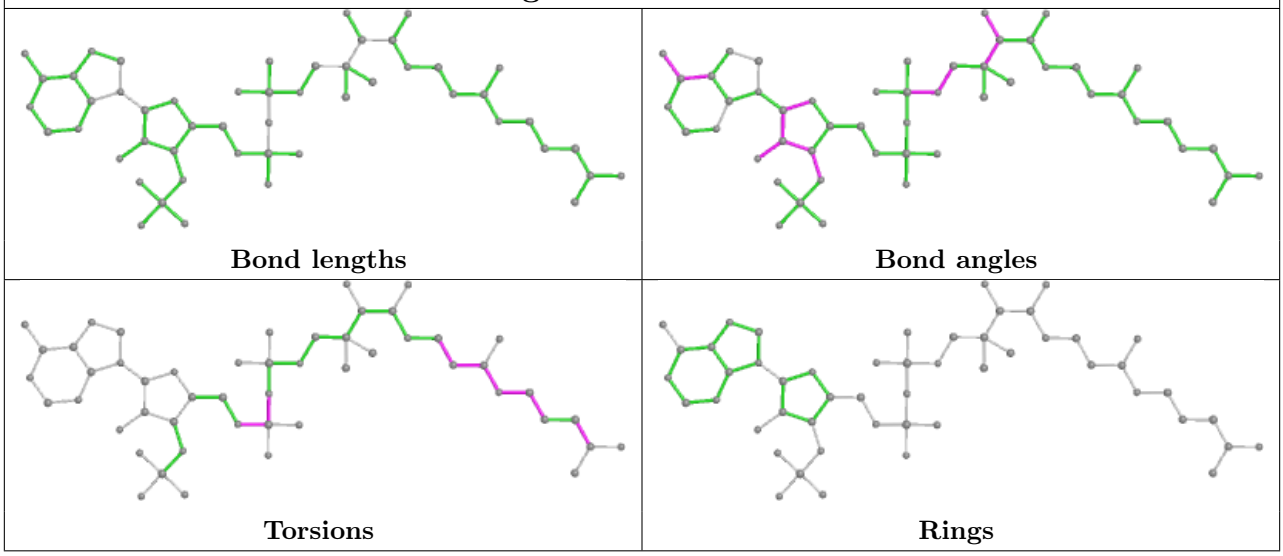
There are no ring outliers.

3 monomers are involved in 3 short contacts:

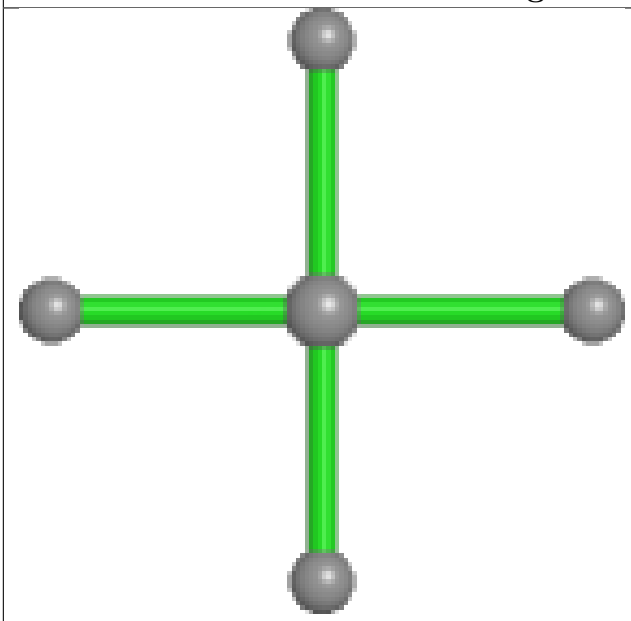
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	903	PO4	1	0
2	B	901	ACO	1	0
2	D	902	ACO	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less then 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

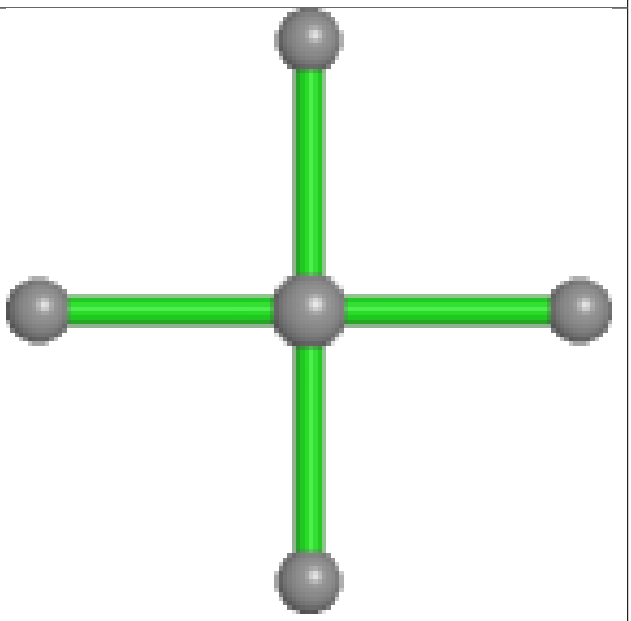
Ligand ACO D 901



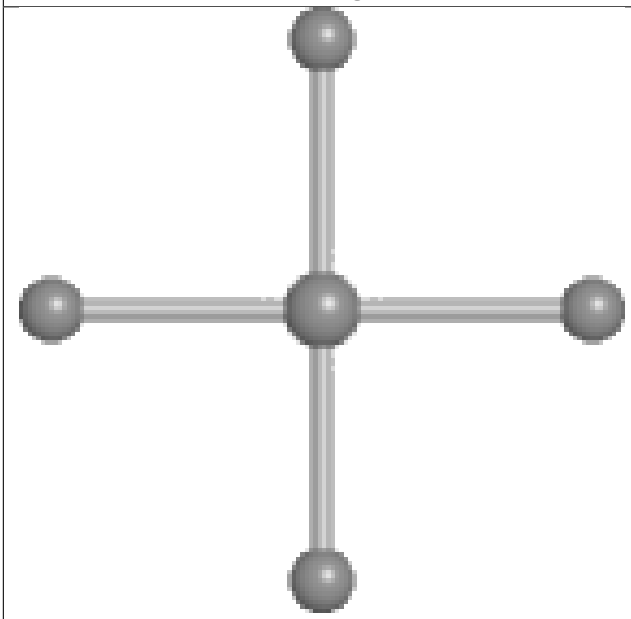
Ligand PO4 B 904



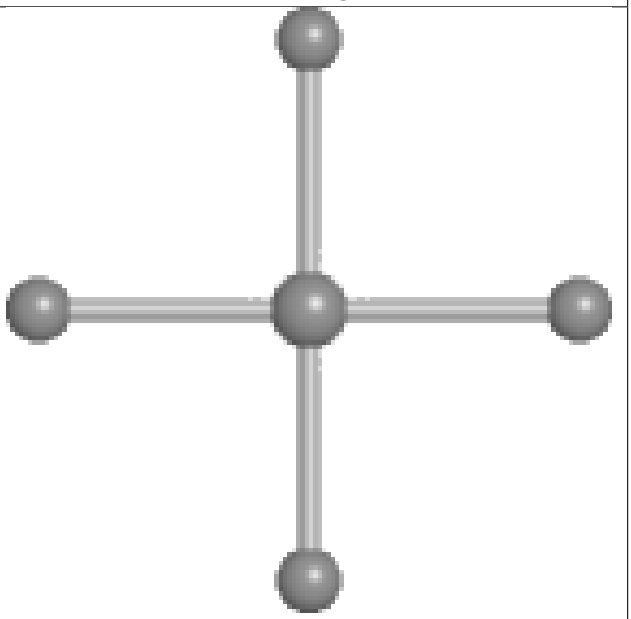
Bond lengths



Bond angles

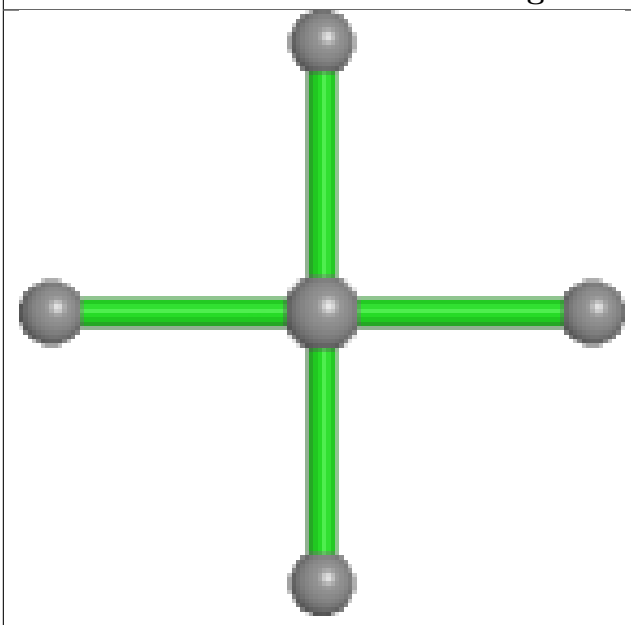


Torsions

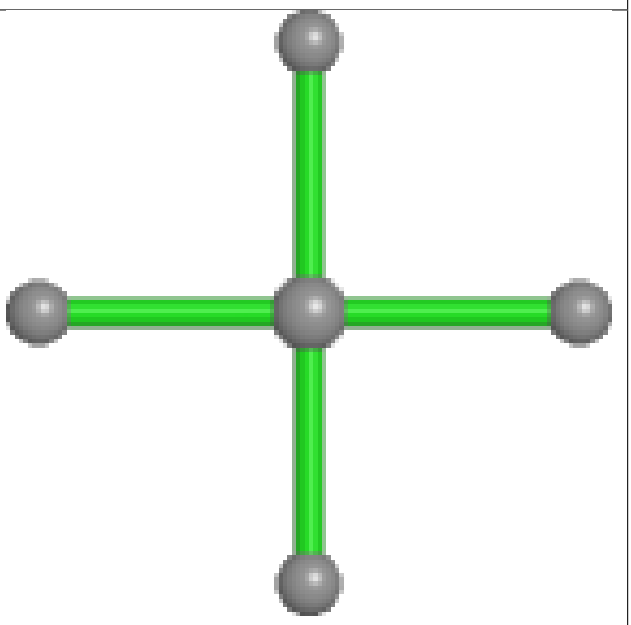


Rings

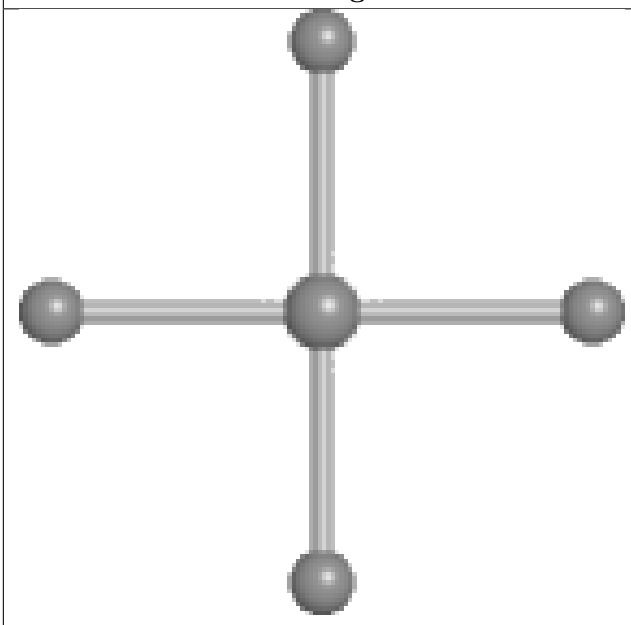
Ligand PO4 C 903



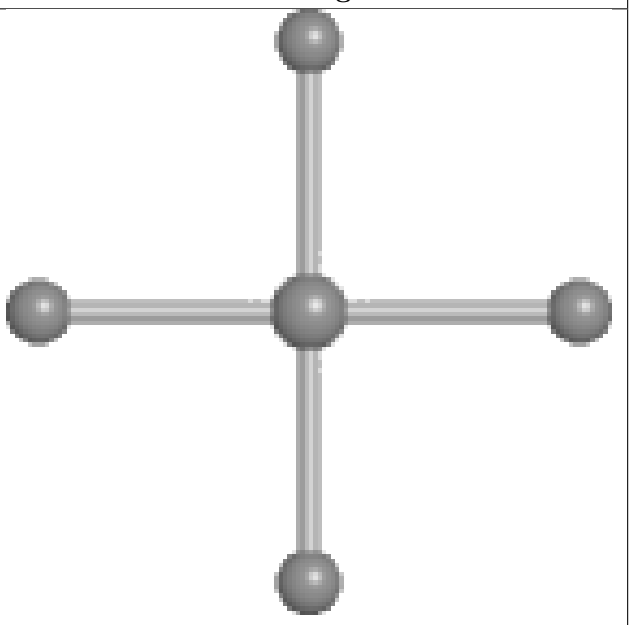
Bond lengths



Bond angles

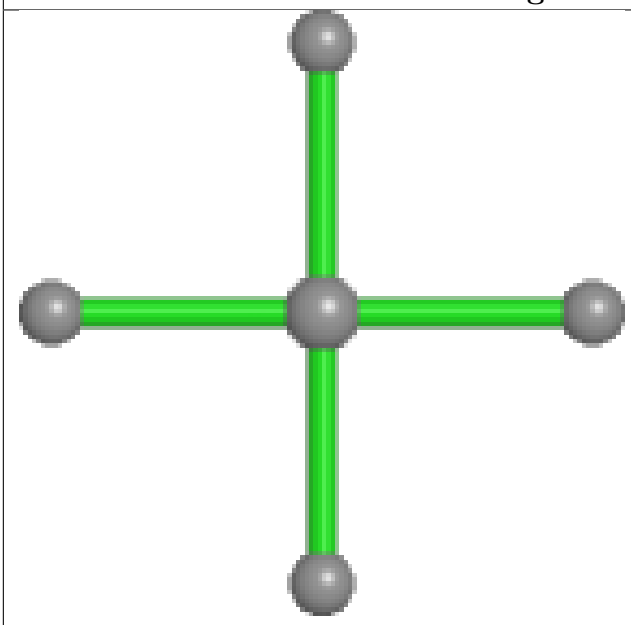


Torsions

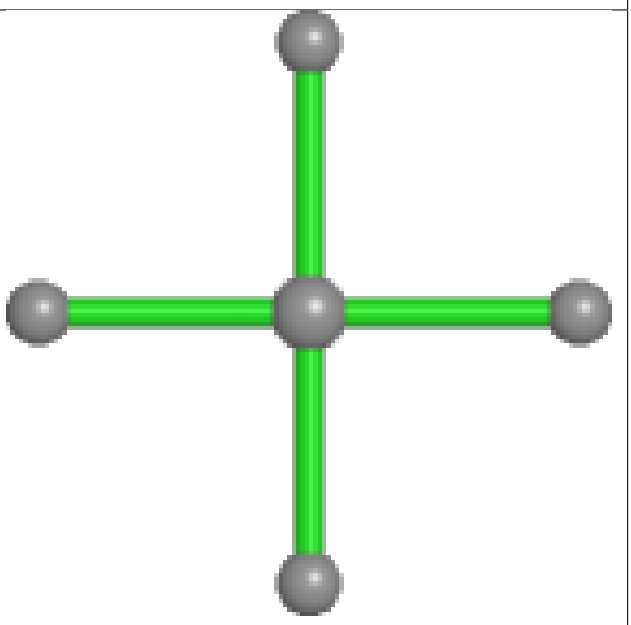


Rings

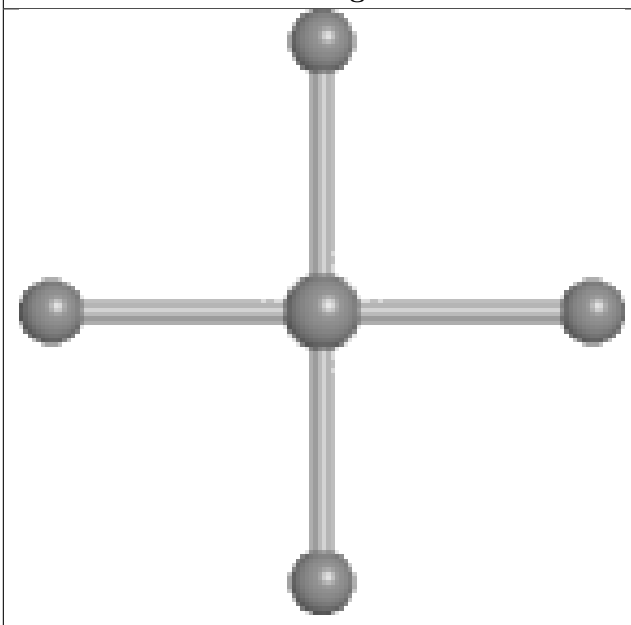
Ligand PO4 C 904



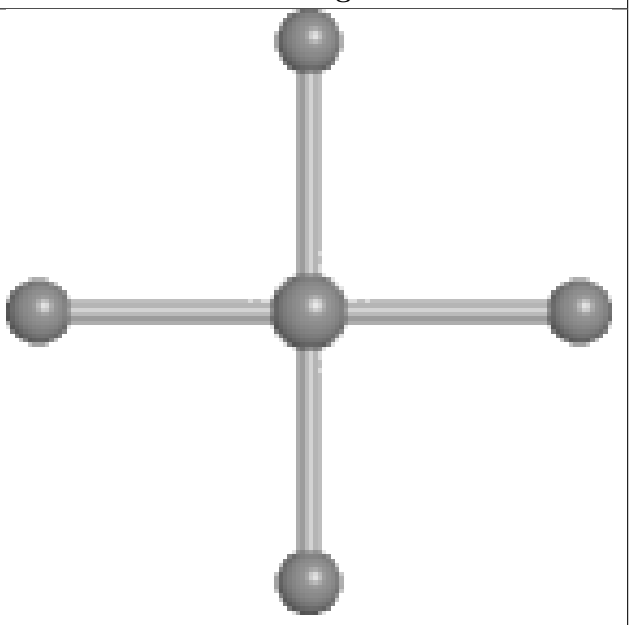
Bond lengths



Bond angles

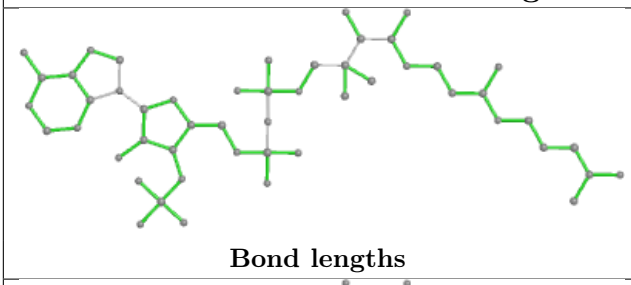


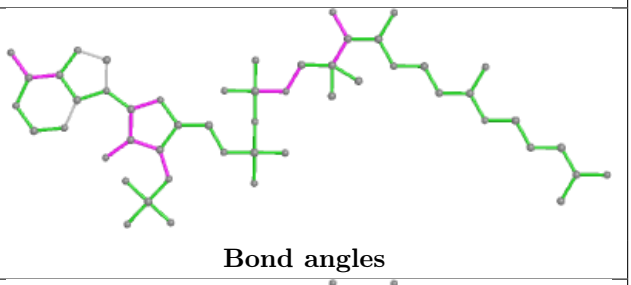
Torsions

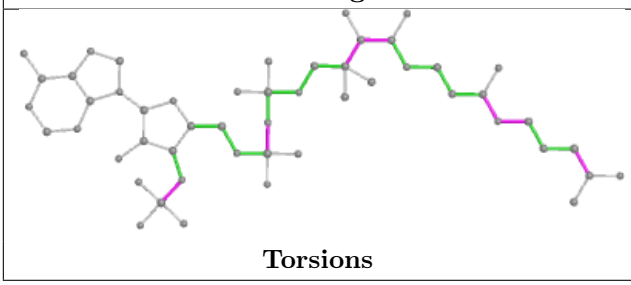


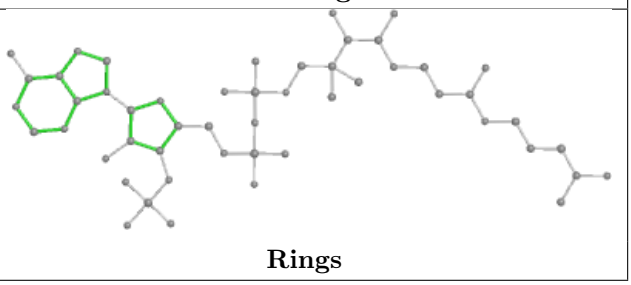
Rings

Ligand ACO B 901

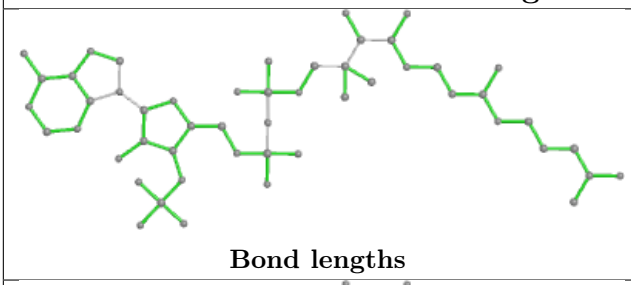

Bond lengths

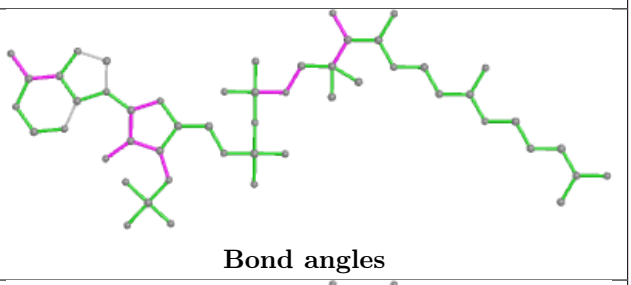

Bond angles

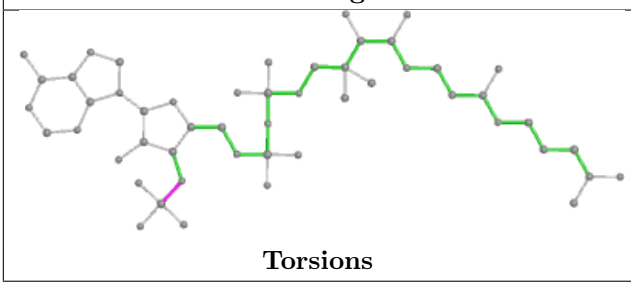

Torsions

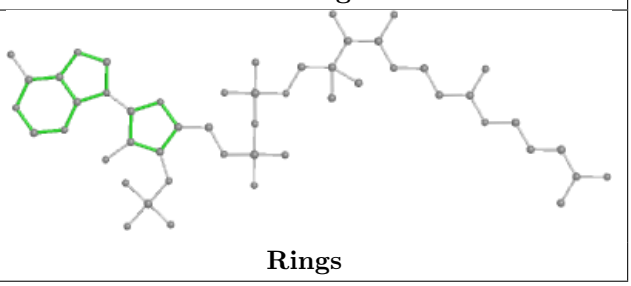

Rings

Ligand ACO C 901

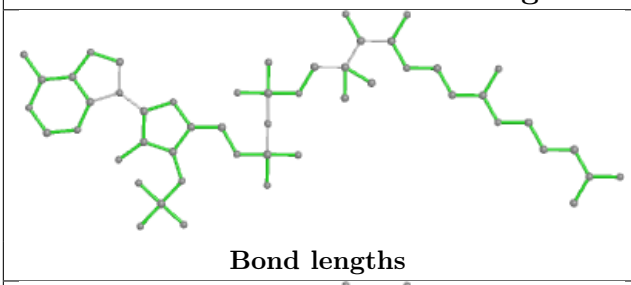

Bond lengths

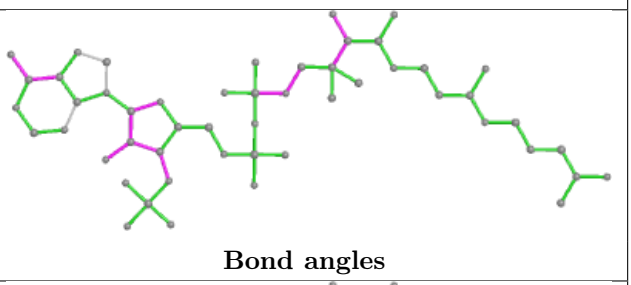

Bond angles

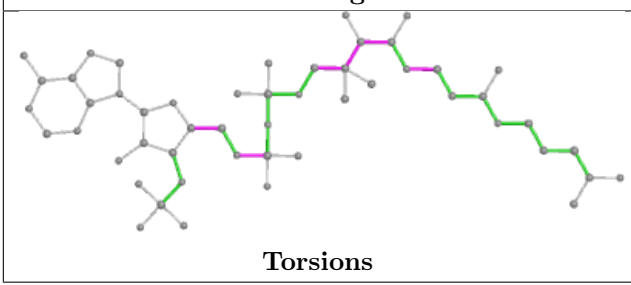

Torsions

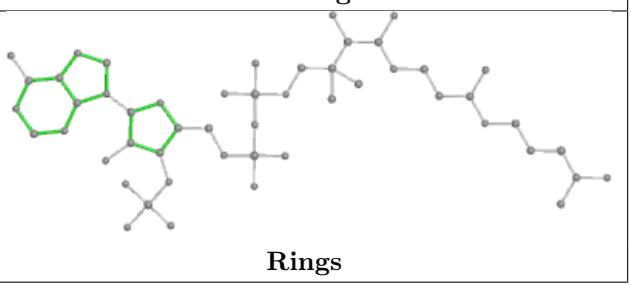

Rings

Ligand ACO A 902

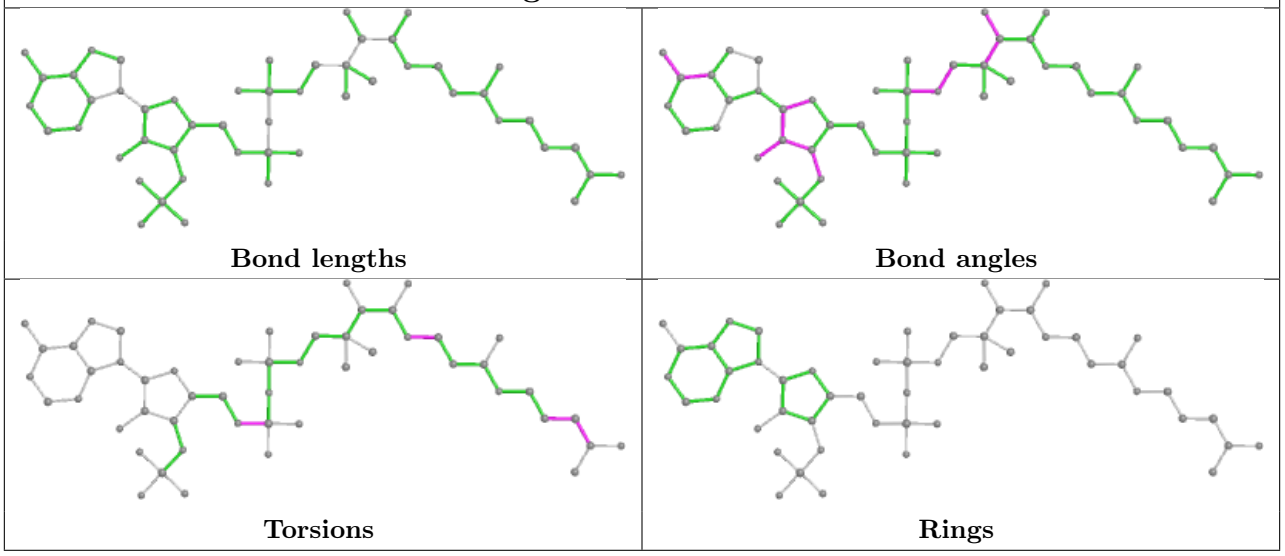

Bond lengths


Bond angles

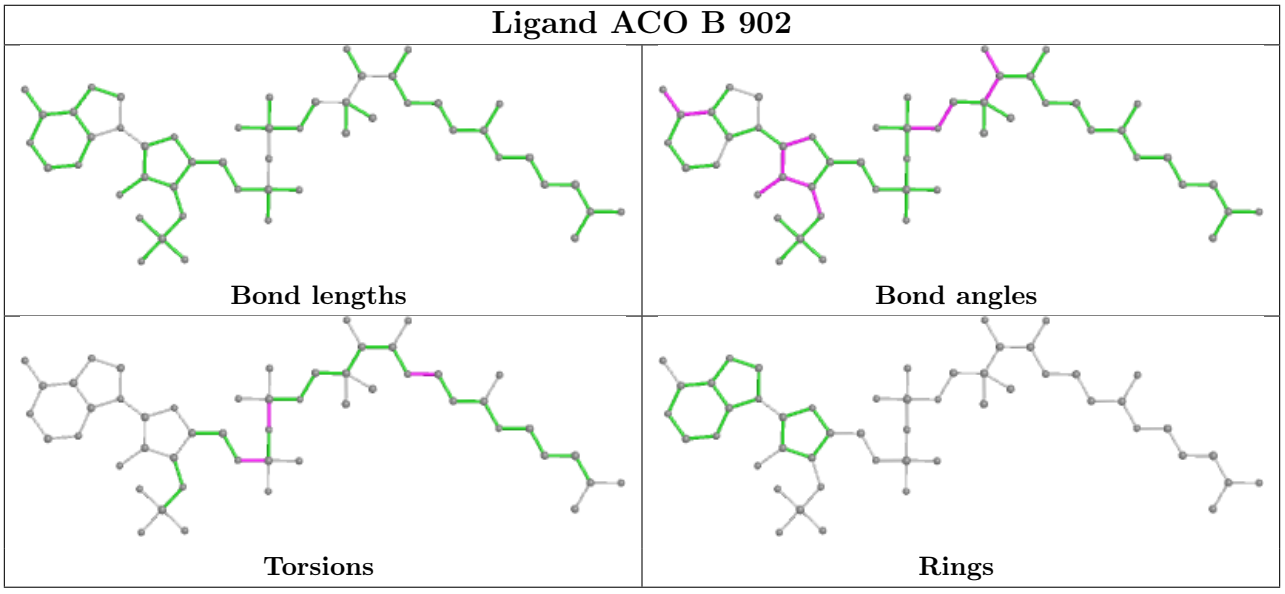

Torsions


Rings

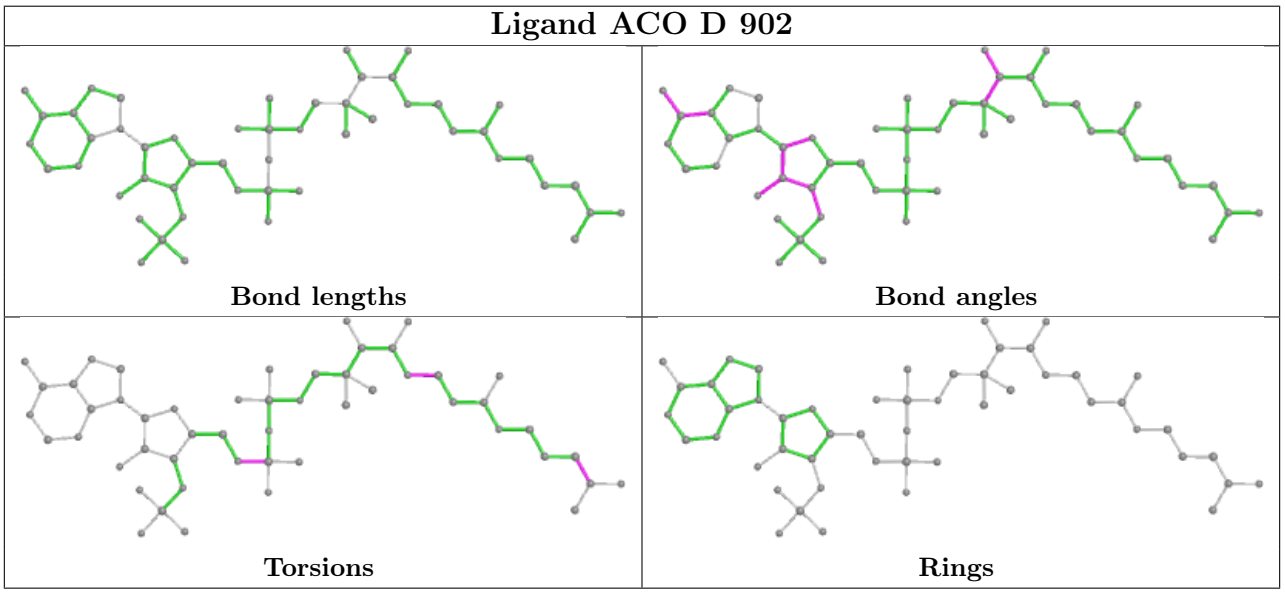
Ligand ACO C 902



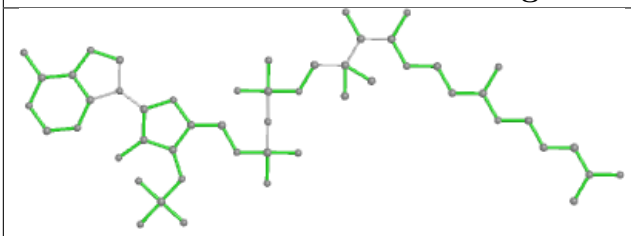
Ligand ACO B 902



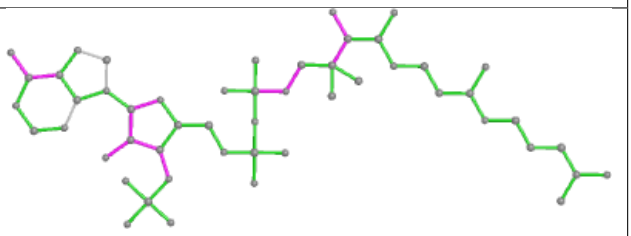
Ligand ACO D 902



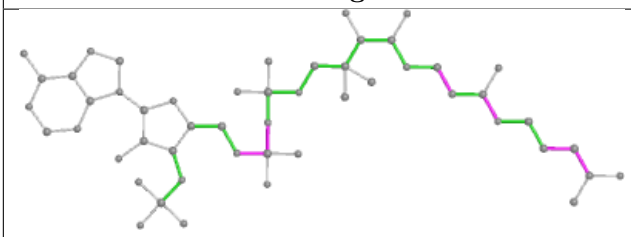
Ligand ACO A 901



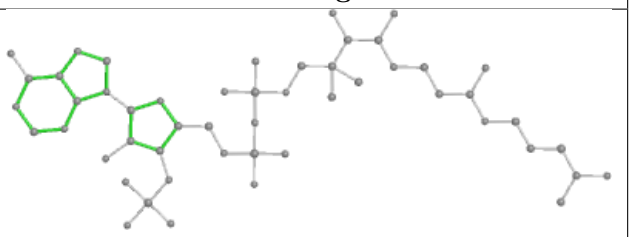
Bond lengths



Bond angles

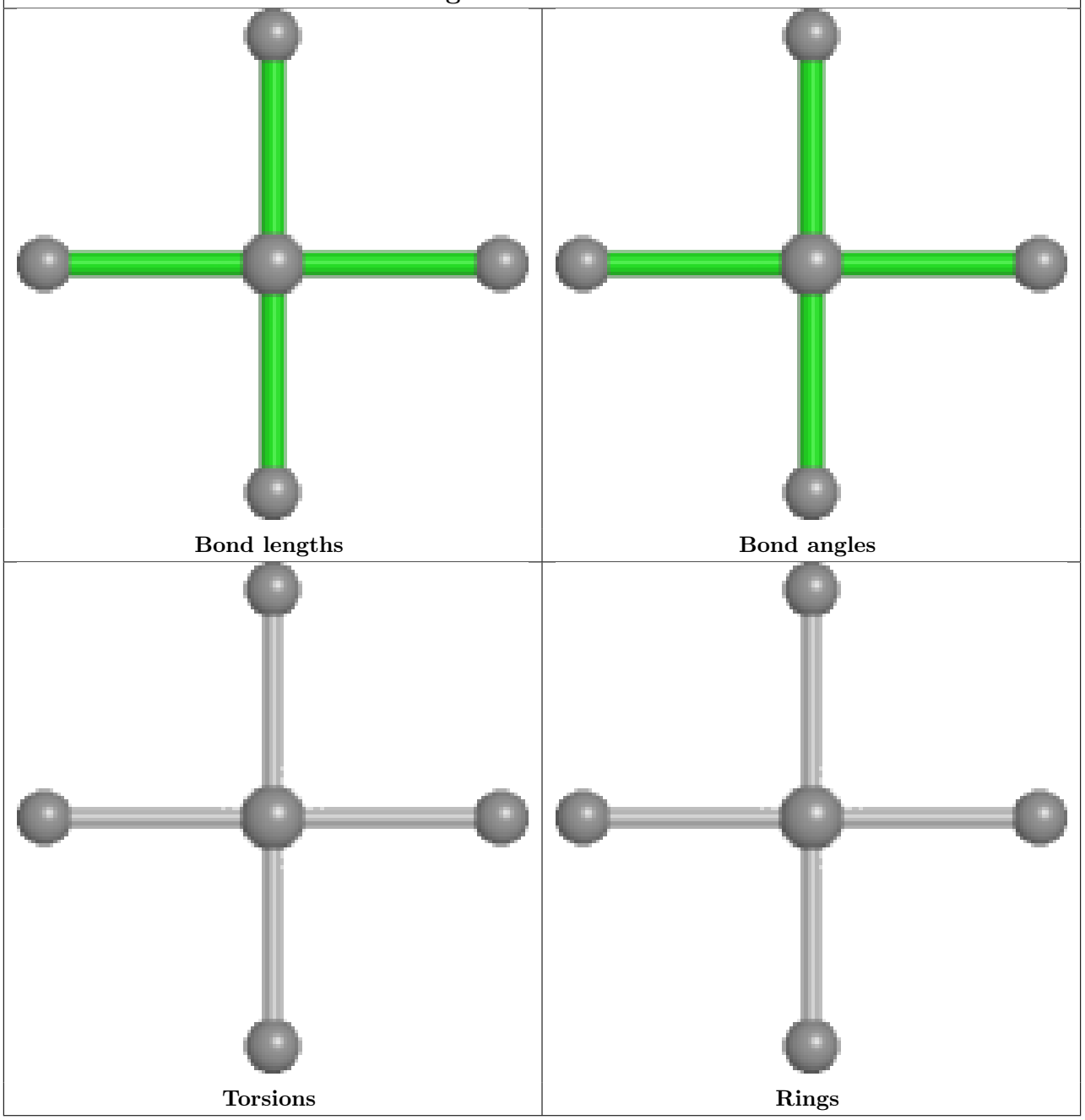


Torsions



Rings

Ligand PO4 B 903



5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

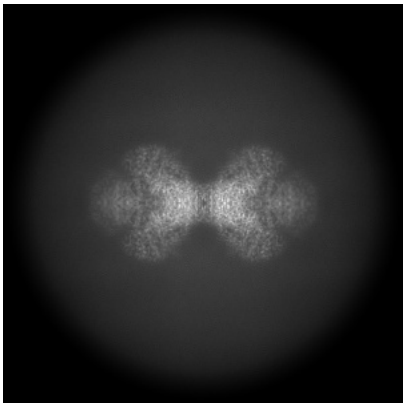
6 Map visualisation

This section contains visualisations of the EMDB entry EMD-60853. These allow visual inspection of the internal detail of the map and identification of artifacts.

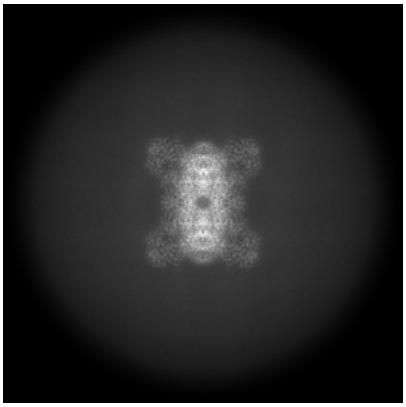
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections

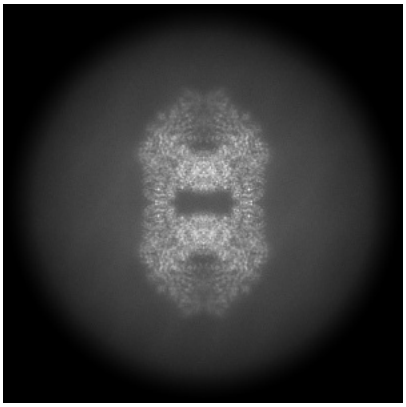
6.1.1 Primary map



X



Y

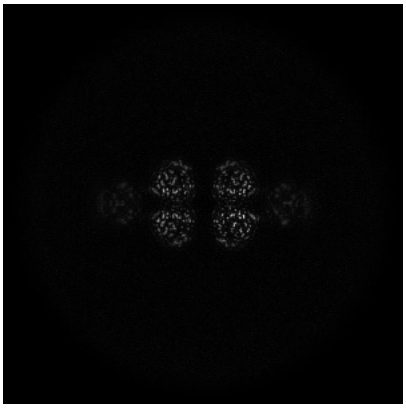


Z

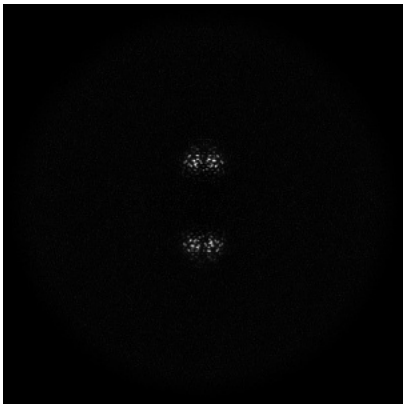
The images above show the map projected in three orthogonal directions.

6.2 Central slices

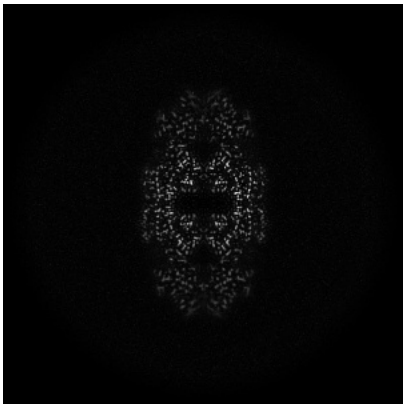
6.2.1 Primary map



X Index: 224



Y Index: 224

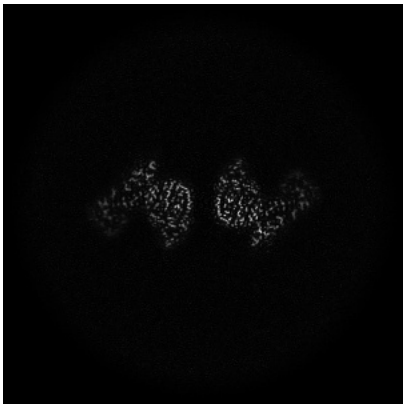


Z Index: 224

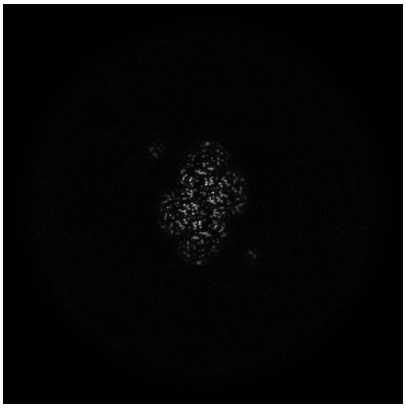
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices (i)

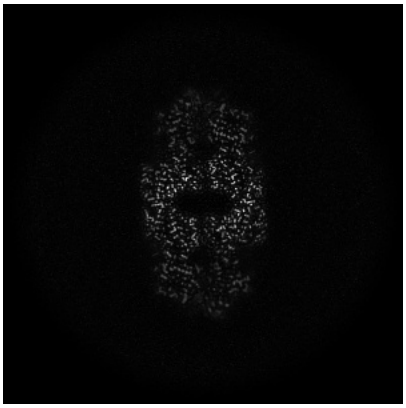
6.3.1 Primary map



X Index: 248



Y Index: 256

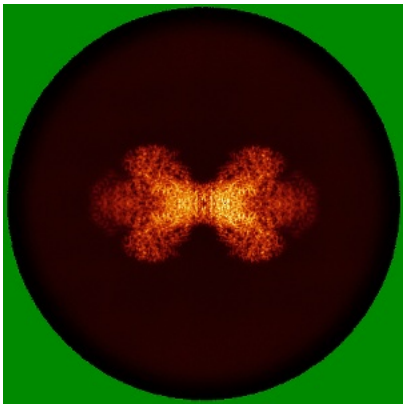


Z Index: 219

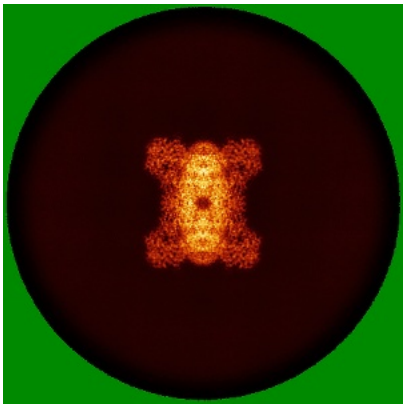
The images above show the largest variance slices of the map in three orthogonal directions.

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

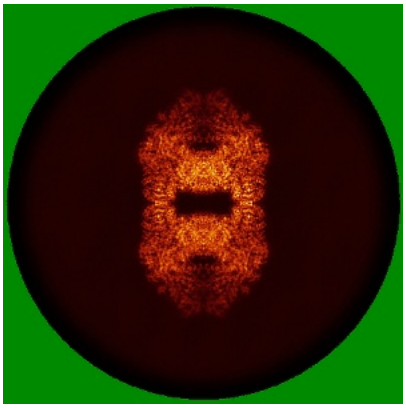
6.4.1 Primary map



X



Y

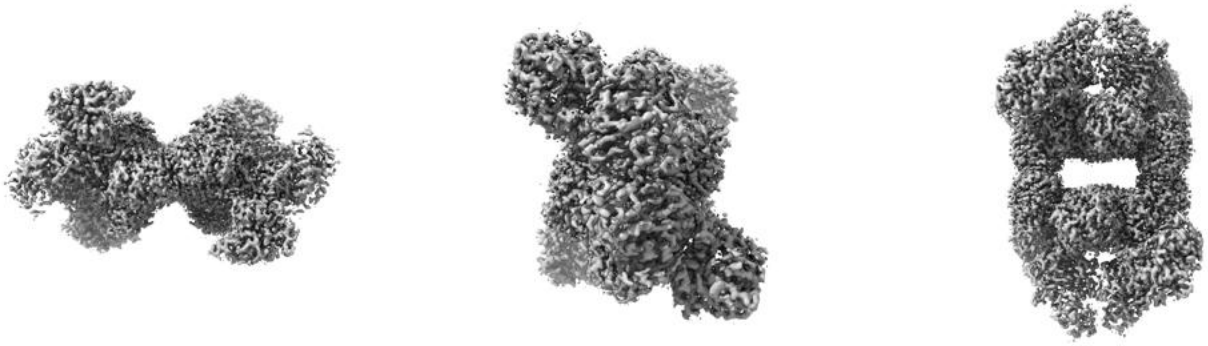


Z

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

6.5 Orthogonal surface views (i)

6.5.1 Primary map



X Y Z

The images above show the 3D surface view of the map at the recommended contour level 0.12. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

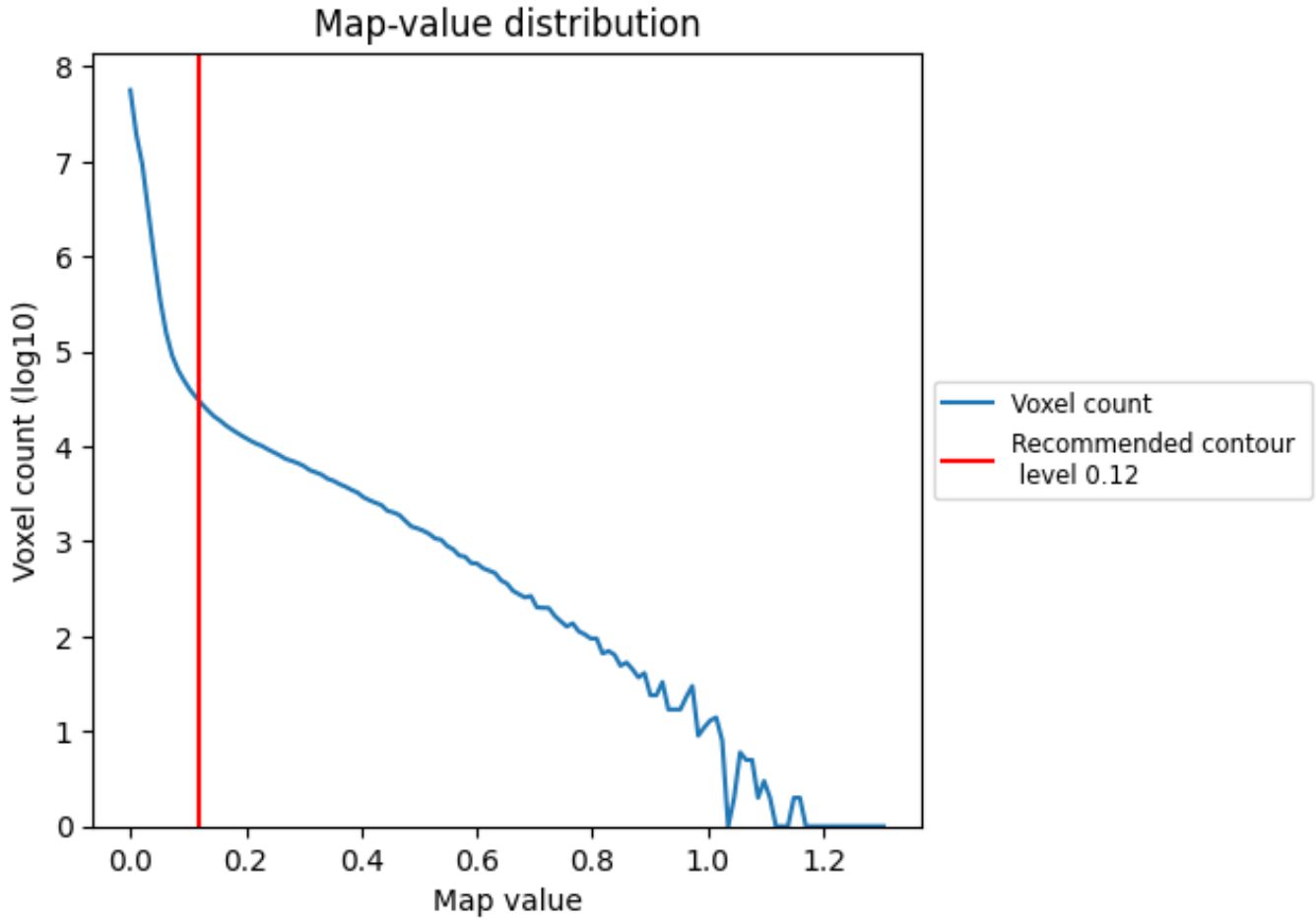
6.6 Mask visualisation (i)

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

7 Map analysis

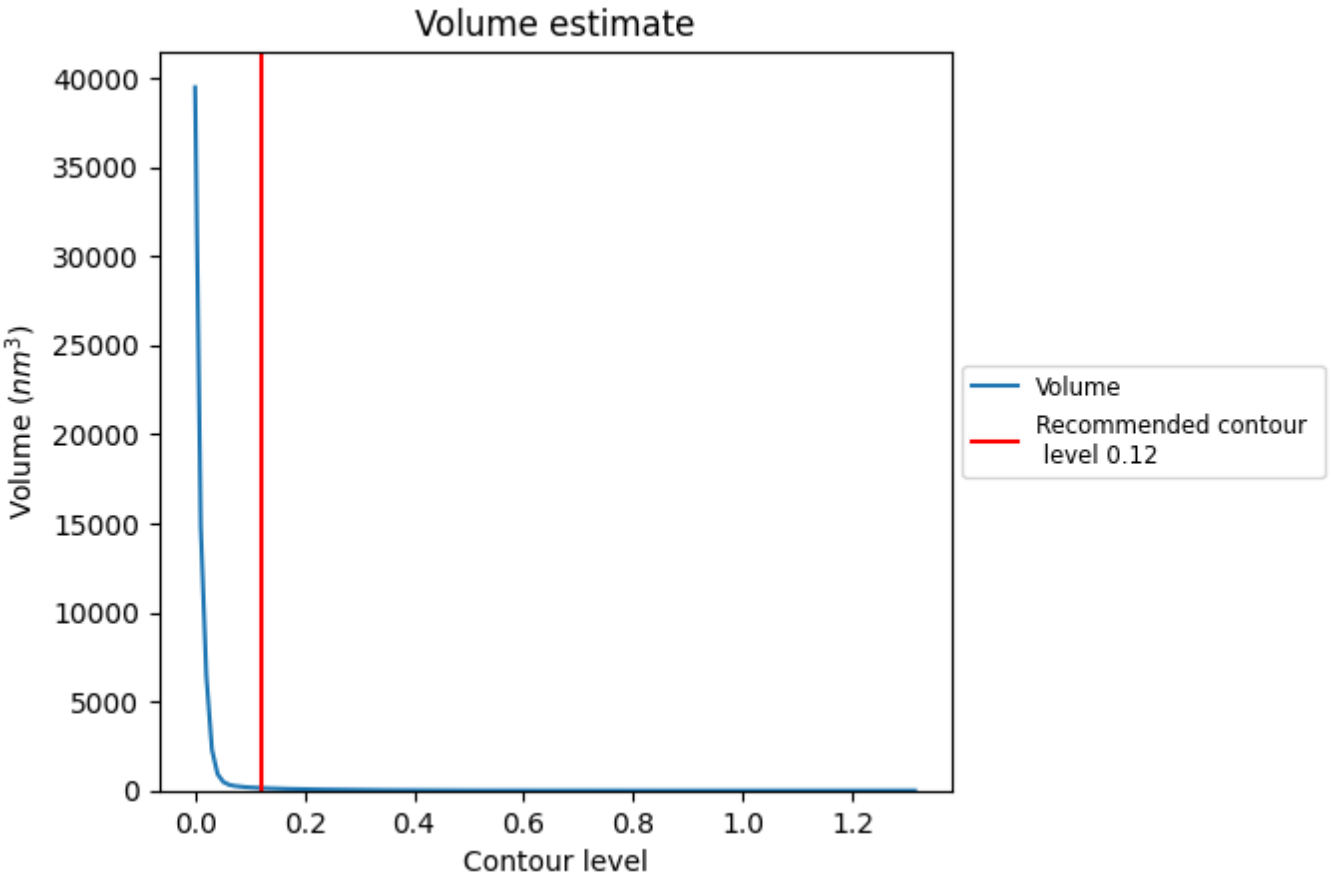
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution



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

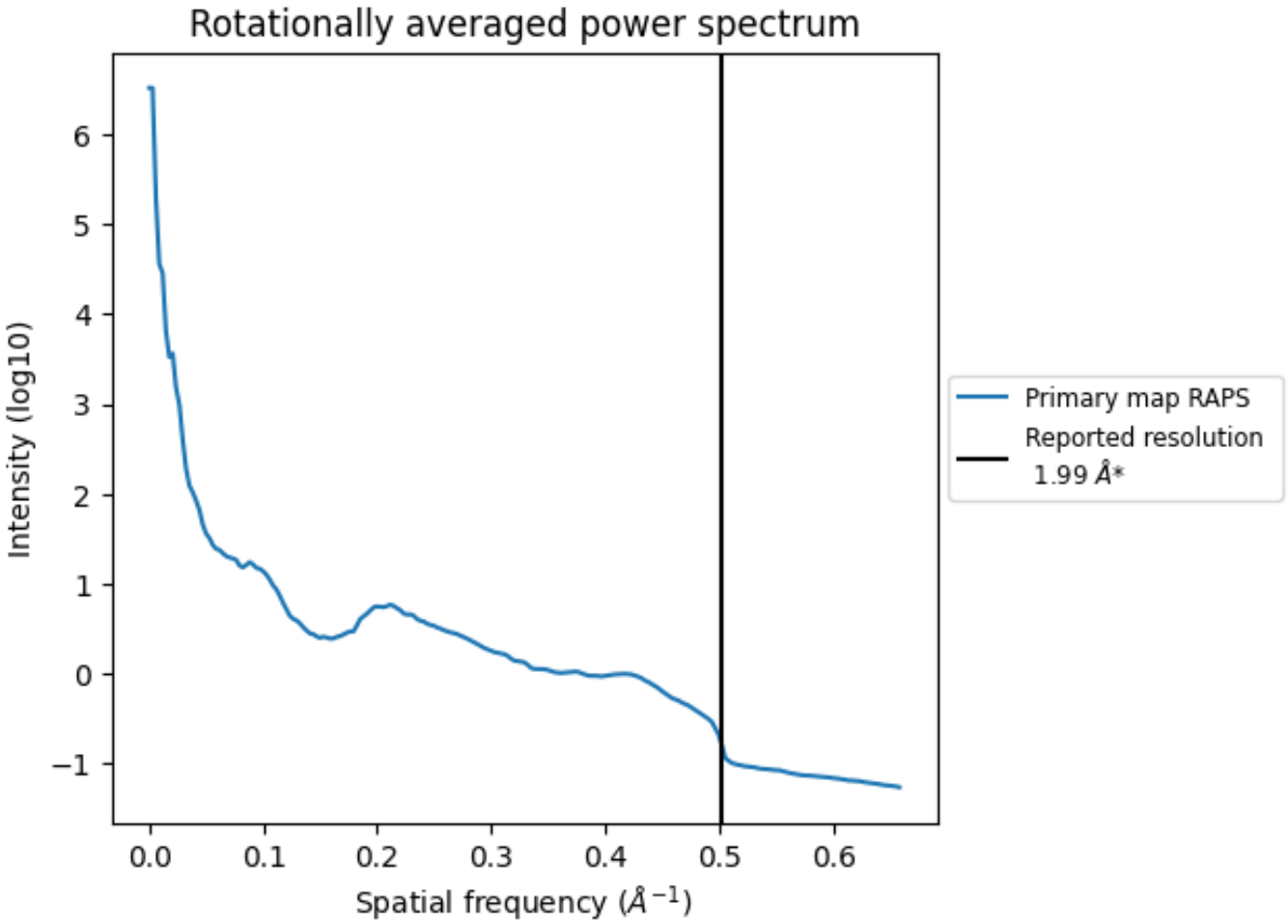
7.2 Volume estimate



The volume at the recommended contour level is 142 nm^3; this corresponds to an approximate mass of 128 kDa.

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

7.3 Rotationally averaged power spectrum (i)



*Reported resolution corresponds to spatial frequency of 0.503 Å^{-1}

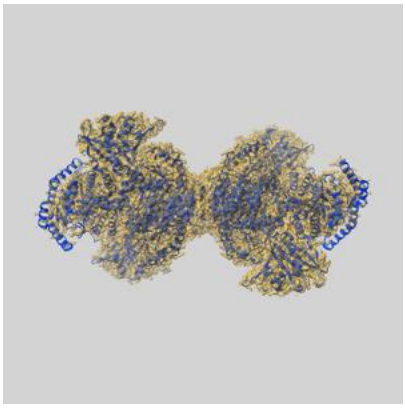
8 Fourier-Shell correlation (i)

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

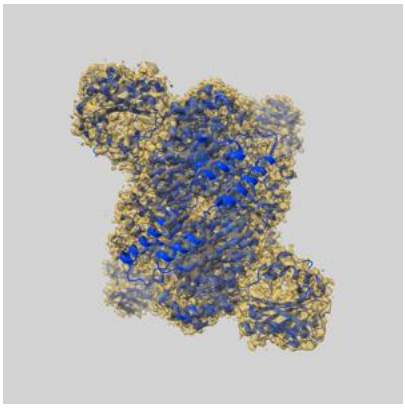
9 Map-model fit

This section contains information regarding the fit between EMDB map EMD-60853 and PDB model 9IT0. Per-residue inclusion information can be found in section 3 on page 7.

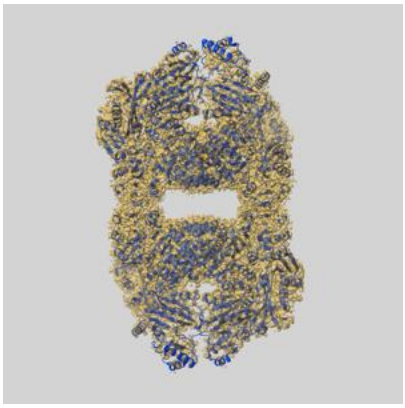
9.1 Map-model overlay



X



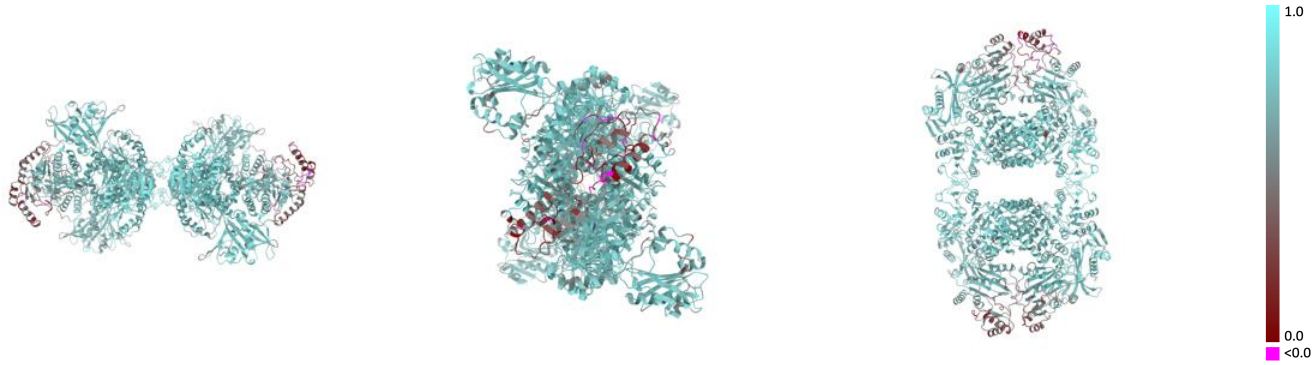
Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.12 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

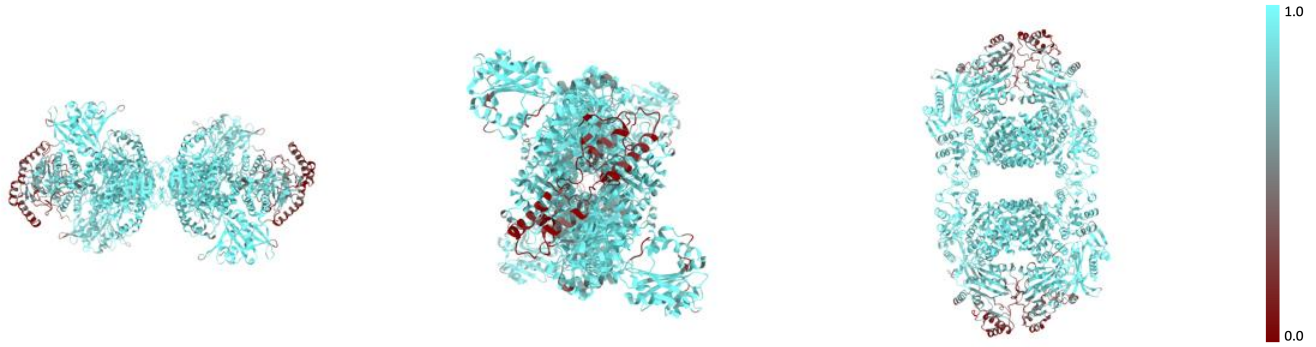
9.2 Q-score mapped to coordinate model (i)



X Y Z

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

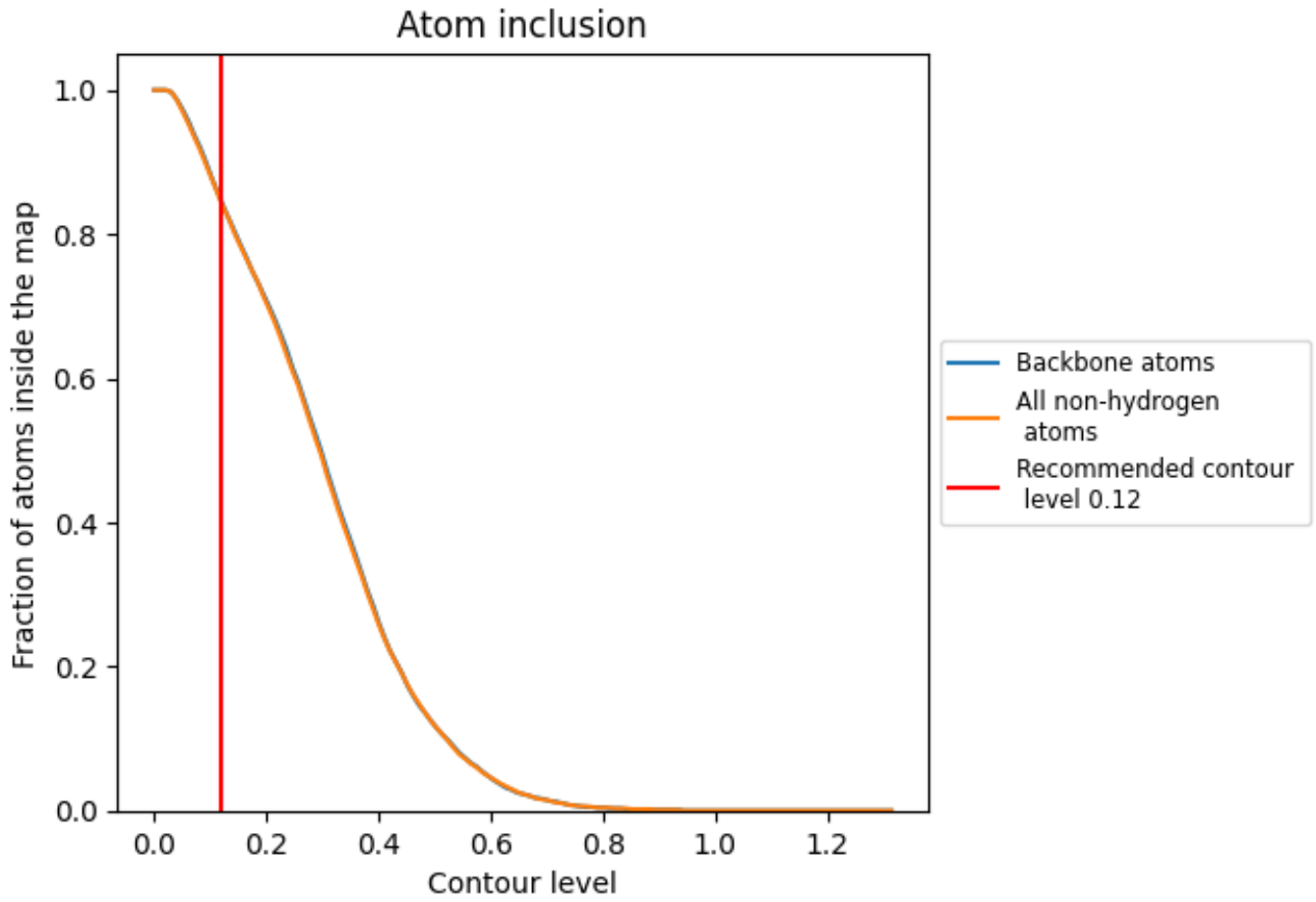
9.3 Atom inclusion mapped to coordinate model (i)



X Y Z

The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (0.12).

9.4 Atom inclusion (i)



At the recommended contour level, 84% of all backbone atoms, 84% 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 (0.12) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.8430	0.7010
A	0.8450	0.7020
B	0.8440	0.6950
C	0.8520	0.7060
D	0.8560	0.7020