Full wwPDB EM Validation Report (i)

Aug 12, 2026 – 09:18 PM JST

PDB ID	:	9UD6 / pdb_00009ud6
EMDB ID	:	EMD-64063
Title	:	Cryo-EM structure of Na+-translocating NADH-ubiquinone oxidoreductase from Vibrio cholerae reduced by NADH, in the absence of Na+, upper state
Authors	:	Ishikawa-Fukuda, M.; Kishikawa, J.; Kato, T.; Murai, M.
Deposited on	:	2025-04-06
Resolution	:	2.65 Å(reported)
Based on initial model	:	7XK3

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 (i) 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 (i)) were used in the production of this report:

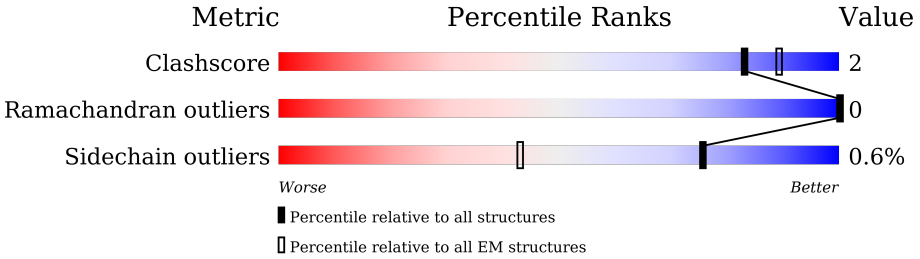
EMDB validation analysis	:	0.0.1.dev133
Mogul	:	1.8.5 (274361), CSD as541be (2020)
MolProbity	:	4-5-2 with Phenix2.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics	:	202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ	:	1.9.13
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.50

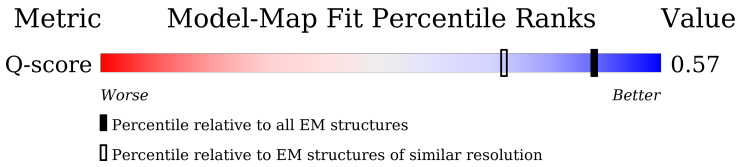
1 Overall quality at a glance (i)

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

The reported resolution of this entry is 2.65 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	9050 (2.15 - 3.15)

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	446	95% 5%
2	B	415	90% • 7%
3	C	257	95% • •
4	D	210	93% • •

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Mol	Chain	Length	Quality of chain
5	E	198	89% 11%
6	F	414	65% 89% 9%

2 Entry composition (i)

There are 13 unique types of molecules in this entry. The entry contains 14743 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 Na(+)-translocating NADH-quinone reductase subunit A.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	A	446	3416	2165	584	650	17	0	0

- Molecule 2 is a protein called Na(+)-translocating NADH-quinone reductase subunit B.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
2	B	387	2964	1960	481	501	22	0	0

- Molecule 3 is a protein called Na(+)-translocating NADH-quinone reductase subunit C.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
3	C	251	1902	1204	327	367	4	0	0

- Molecule 4 is a protein called Na(+)-translocating NADH-quinone reductase subunit D.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
4	D	204	1562	1037	247	268	10	0	0

- Molecule 5 is a protein called Na(+)-translocating NADH-quinone reductase subunit E.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
5	E	198	1511	1013	230	257	11	0	0

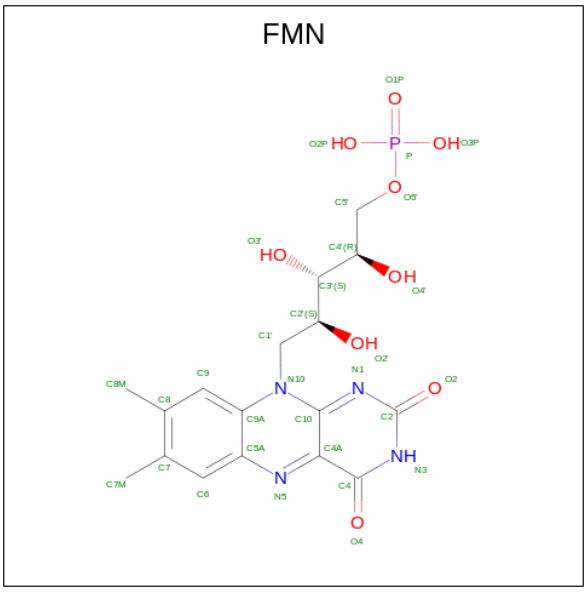
- Molecule 6 is a protein called Na(+)-translocating NADH-quinone reductase subunit F.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
6	F	408	3165	2027	519	595	24	0	0

There are 6 discrepancies between the modelled and reference sequences:

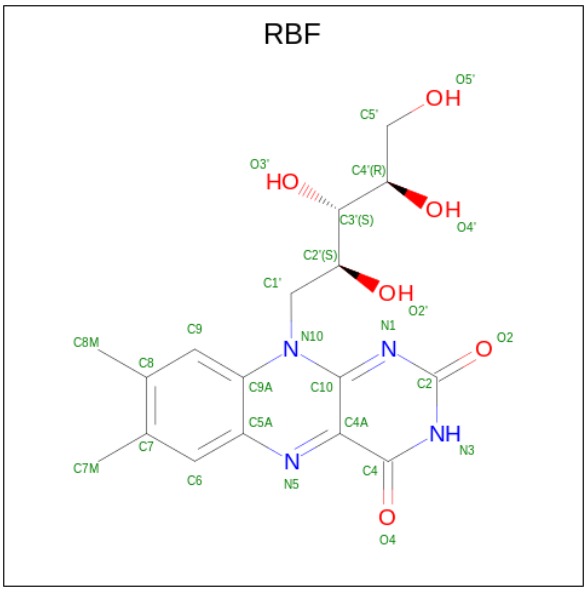
Chain	Residue	Modelled	Actual	Comment	Reference
F	409	HIS	-	expression tag	UNP A5F5Y4
F	410	HIS	-	expression tag	UNP A5F5Y4
F	411	HIS	-	expression tag	UNP A5F5Y4
F	412	HIS	-	expression tag	UNP A5F5Y4
F	413	HIS	-	expression tag	UNP A5F5Y4
F	414	HIS	-	expression tag	UNP A5F5Y4

- Molecule 7 is FLAVIN MONONUCLEOTIDE (CCD ID: FMN) (formula: $C_{17}H_{21}N_4O_9P$).



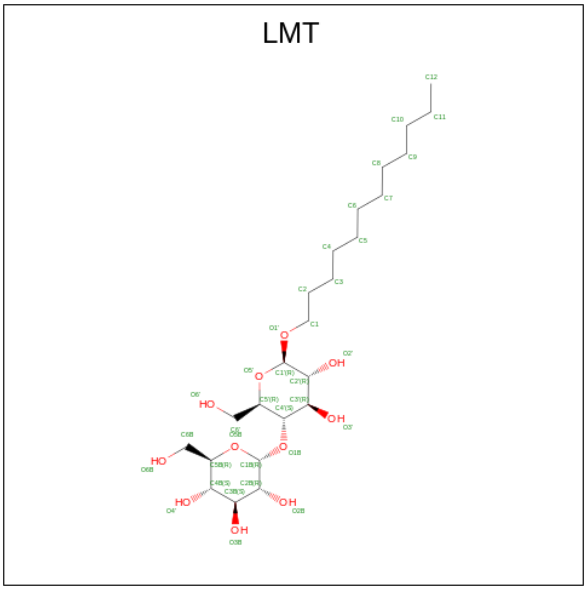
Mol	Chain	Residues	Atoms					AltConf
			Total	C	N	O	P	
7	B	1	31	17	4	9	1	0
			Total	C	N	O	P	
7	C	1	31	17	4	9	1	0

- Molecule 8 is RIBOFLAVIN (CCD ID: RBF) (formula: $C_{17}H_{20}N_4O_6$).



Mol	Chain	Residues	Atoms				AltConf
			Total	C	N	O	
8	B	1	27	17	4	6	0

- Molecule 9 is DODECYL-BETA-D-MALTOSIDE (CCD ID: LMT) (formula: $C_{24}H_{46}O_{11}$).

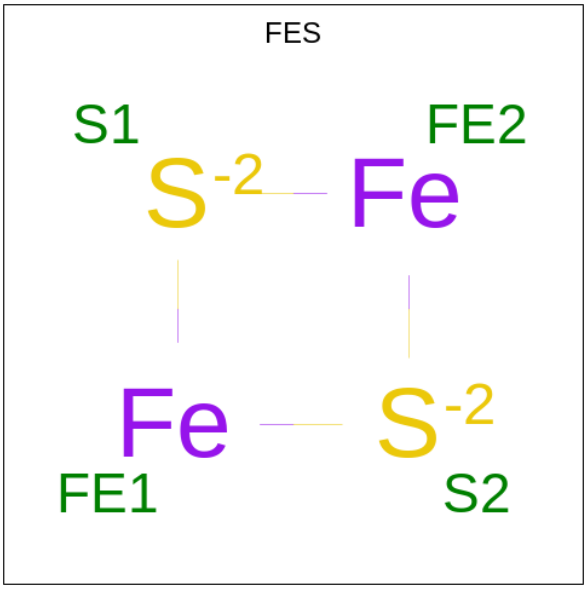


Mol	Chain	Residues	Atoms				AltConf
			Total	C	O		
9	B	1	35	24	11		0
			Total	C	O		
9	D	1	35	24	11		0

- Molecule 10 is CALCIUM ION (CCD ID: CA) (formula: Ca).

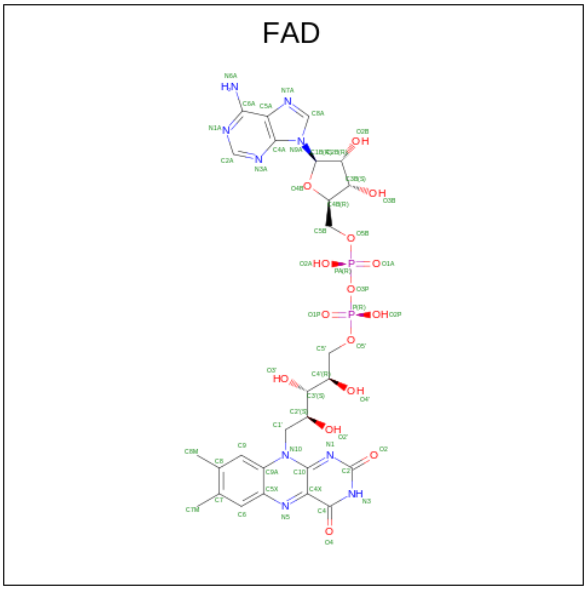
Mol	Chain	Residues	Atoms		AltConf
			Total	Ca	
10	C	1	1	1	0

- Molecule 11 is FE2/S2 (INORGANIC) CLUSTER (CCD ID: FES) (formula: Fe_2S_2).



Mol	Chain	Residues	Atoms			AltConf
			Total	Fe	S	
11	E	1	4	2	2	0
11	F	1	4	2	2	0

- Molecule 12 is FLAVIN-ADENINE DINUCLEOTIDE (CCD ID: FAD) (formula: $C_{27}H_{33}N_9O_{15}P_2$).



Mol	Chain	Residues	Atoms					AltConf
			Total	C	N	O	P	
12	F	1	53	27	9	15	2	0

- Molecule 13 is water.

Mol	Chain	Residues	Atoms		AltConf
			Total	O	
13	D	2	2	2	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: Na(+)-translocating NADH-quinone reductase subunit A

Chain A: 95% 5%

M1 G14 M39 R40 I56 E59 V72 V94 R115 T133 R134 S137 S146 L159 T185 T186 K193 T196 M224 P229 M339 D396 S404 K444 E445 G446

• Molecule 2: Na(+)-translocating NADH-quinone reductase subunit B

Chain B: 90% 7%

MET GLY LEU LYS LYS PHE LEU GLU GLU ASP ILE GLU HIS HIS PHE GLU PRO GLY GLY LYS HIS GLU LYS TRP PHE ALA LEU Y27 E28 A29 A30 A31 T32 R43 D49 D52 W100 W103 V136 R151 I183 V193 T236 A237 L333 M382 V401

R411 Y412 G413 LYS GLN

• Molecule 3: Na(+)-translocating NADH-quinone reductase subunit C

Chain C: 95%

MET ALA SER ASN ASN D6 V31 D131 L145 E197 A210 P211 Q212 D219 L226 L256 ASN

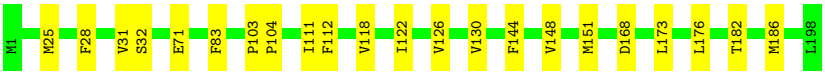
• Molecule 4: Na(+)-translocating NADH-quinone reductase subunit D

Chain D: 93%

MET SER SER ALA LYS GLU L7 L38 I72 Q75 M115 I128 P129 L145 M146 M181 E210

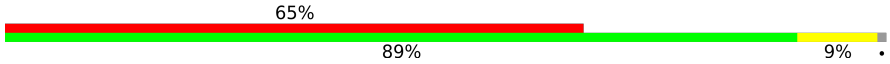
• Molecule 5: Na(+)-translocating NADH-quinone reductase subunit E

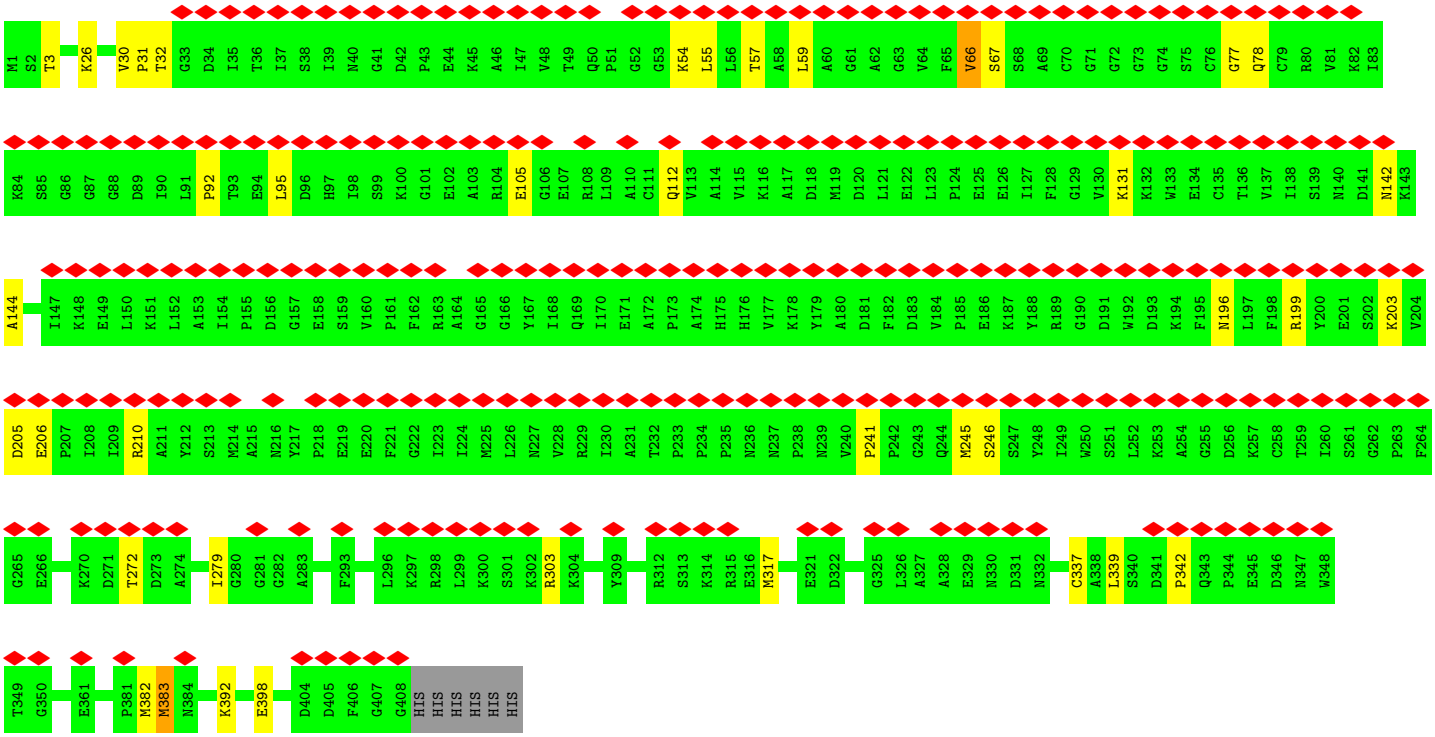
Chain E: 89% 11%



• Molecule 6: Na(+)-translocating NADH-quinone reductase subunit F

Chain F:





4 Experimental information (i)

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	1244065	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 ($e^-/Å^2$)	65	Depositor
Minimum defocus (nm)	800	Depositor
Maximum defocus (nm)	1800	Depositor
Magnification	81000	Depositor
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	3.126	Depositor
Minimum map value	-2.183	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.059	Depositor
Recommended contour level	0.195	Depositor
Map size (Å)	281.6, 281.6, 281.6	wwPDB
Map dimensions	320, 320, 320	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.88, 0.88, 0.88	Depositor

5 Model quality (i)

5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: FMN, LMT, FES, RBF, CA, FAD

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	0.16	0/3480	0.30	0/4715
2	B	0.17	0/3053	0.32	0/4160
3	C	0.15	0/1934	0.28	0/2610
4	D	0.16	0/1594	0.30	0/2164
5	E	0.16	0/1544	0.31	0/2094
6	F	0.11	0/3243	0.32	2/4389 (0.0%)
All	All	0.15	0/14848	0.31	2/20132 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
6	F	66	VAL	CA-C-N	6.26	132.97	121.70
6	F	66	VAL	C-N-CA	6.26	132.97	121.70

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	3416	0	3463	13	0
2	B	2964	0	2952	8	0
3	C	1902	0	1911	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	D	1562	0	1648	6	0
5	E	1511	0	1586	14	0
6	F	3165	0	3103	27	0
7	B	31	0	19	2	0
7	C	31	0	19	0	0
8	B	27	0	20	0	0
9	B	35	0	46	0	0
9	D	35	0	46	4	0
10	C	1	0	0	0	0
11	E	4	0	0	0	0
11	F	4	0	0	0	0
12	F	53	0	31	4	0
13	D	2	0	0	0	0
All	All	14743	0	14844	72	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 2.

All (72) 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)
6:F:206:GLU:N	6:F:206:GLU:OE1	2.18	0.76
9:D:301:LMT:H2O1	9:D:301:LMT:H6'	0.75	0.73
7:B:501:FMN:O2P	7:B:501:FMN:O4'	2.10	0.70
6:F:105:GLU:N	6:F:105:GLU:OE1	2.26	0.69
6:F:196:ASN:O	6:F:199:ARG:NH1	2.26	0.68
1:A:193:LYS:NZ	1:A:196:THR:O	2.28	0.66
6:F:31:PRO:O	6:F:32:THR:OG1	2.09	0.64
2:B:151:ARG:NH2	5:E:168:ASP:O	2.31	0.64
5:E:31:VAL:HG12	5:E:31:VAL:O	2.00	0.61
5:E:28:PHE:O	5:E:32:SER:OG	2.21	0.58
6:F:392:LYS:NZ	6:F:398:GLU:OE2	2.22	0.57
6:F:246:SER:N	12:F:502:FAD:O2P	2.38	0.56
9:D:301:LMT:O6'	9:D:301:LMT:O2B	2.15	0.55
4:D:181:MET:HE3	9:D:301:LMT:H81	1.89	0.55
6:F:142:ASN:ND2	6:F:144:ALA:O	2.35	0.55
6:F:112:GLN:N	6:F:112:GLN:OE1	2.41	0.54
6:F:66:VAL:N	6:F:67:SER:O	2.41	0.54
6:F:66:VAL:HB	6:F:67:SER:HB3	1.89	0.53
1:A:72:VAL:HG21	1:A:94:VAL:HG22	1.92	0.52
2:B:193:VAL:HG22	5:E:173:LEU:HD13	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:F:54:LYS:O	6:F:57:THR:OG1	2.27	0.50
3:C:6:ASP:N	3:C:6:ASP:OD1	2.44	0.49
6:F:241:PRO:O	12:F:502:FAD:O2B	2.16	0.49
5:E:25:MET:HA	5:E:25:MET:HE2	1.95	0.49
6:F:26:LYS:O	6:F:30:VAL:HG22	2.12	0.49
1:A:396:ASP:OD2	1:A:404:SER:OG	2.31	0.49
4:D:181:MET:HE2	9:D:301:LMT:H101	1.96	0.48
2:B:52:ASP:OD1	2:B:52:ASP:N	2.45	0.48
1:A:339:MET:HA	1:A:339:MET:HE2	1.96	0.48
6:F:203:LYS:NZ	6:F:205:ASP:OD1	2.44	0.47
6:F:77:GLY:HA2	6:F:78:GLN:HB2	1.96	0.47
6:F:245:MET:N	12:F:502:FAD:O2P	2.46	0.47
2:B:136:VAL:HG13	2:B:183:ILE:HD13	1.97	0.45
3:C:31:VAL:HG21	6:F:3:THR:HG23	1.98	0.45
5:E:122:ILE:O	5:E:126:VAL:HG23	2.16	0.45
1:A:133:THR:O	1:A:137:SER:N	2.49	0.45
4:D:75:GLN:OE1	4:D:115:MET:HE1	2.17	0.45
6:F:92:PRO:HA	6:F:95:LEU:HD23	1.98	0.45
3:C:145:LEU:HD13	3:C:226:LEU:HD13	1.97	0.45
2:B:237:ALA:HB3	2:B:333:LEU:HD22	1.99	0.45
1:A:229:PRO:O	2:B:411:ARG:NH1	2.50	0.44
5:E:71:GLU:HA	5:E:71:GLU:OE2	2.17	0.44
4:D:128:ILE:HB	4:D:129:PRO:HD3	1.98	0.44
6:F:55:LEU:HD13	6:F:59:LEU:HD23	1.99	0.44
1:A:56:ILE:HD11	1:A:59:GLU:HG3	1.99	0.44
2:B:43:ARG:HH11	2:B:43:ARG:HG2	1.83	0.44
5:E:144:PHE:O	5:E:148:VAL:HG23	2.18	0.44
5:E:103:PRO:N	5:E:104:PRO:CD	2.81	0.43
4:D:72:ILE:CD1	4:D:115:MET:HE2	2.48	0.43
5:E:83:PHE:CZ	5:E:130:VAL:HG21	2.54	0.43
2:B:100:TRP:HA	2:B:103:TRP:NE1	2.34	0.42
5:E:111:ILE:HG23	5:E:112:PHE:CD2	2.53	0.42
7:B:501:FMN:O4'	7:B:501:FMN:P	2.78	0.42
4:D:38:LEU:HD22	4:D:146:MET:HE3	2.01	0.42
6:F:66:VAL:N	6:F:67:SER:C	2.78	0.42
6:F:272:THR:OG1	6:F:303:ARG:NH2	2.53	0.42
5:E:182:THR:O	5:E:186:MET:HG3	2.19	0.41
6:F:339:LEU:HD13	6:F:342:PRO:HB3	2.01	0.41
1:A:39:MET:O	1:A:40:ARG:NH1	2.53	0.41
1:A:14:GLY:O	1:A:159:LEU:HD22	2.20	0.41
6:F:95:LEU:N	6:F:95:LEU:HD22	2.35	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:F:210:ARG:NE	12:F:502:FAD:O1P	2.52	0.41
6:F:317:MET:HE1	6:F:337:CYS:SG	2.61	0.41
6:F:382:MET:C	6:F:382:MET:SD	3.03	0.41
1:A:134:ARG:HG3	1:A:224:MET:SD	2.61	0.41
6:F:279:ILE:HG22	6:F:383:MET:HE3	2.02	0.41
3:C:210:ALA:N	3:C:219:ASP:OD1	2.39	0.41
1:A:115:ARG:NE	1:A:186:THR:OG1	2.53	0.41
5:E:118:VAL:O	5:E:118:VAL:CG1	2.69	0.41
1:A:186:THR:HG22	1:A:186:THR:O	2.21	0.41
1:A:146:SER:O	1:A:185:THR:O	2.40	0.40
5:E:151:MET:HE2	5:E:151:MET:HB3	1.99	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	444/446 (100%)	425 (96%)	19 (4%)	0	100	100
2	B	385/415 (93%)	376 (98%)	9 (2%)	0	100	100
3	C	249/257 (97%)	239 (96%)	10 (4%)	0	100	100
4	D	202/210 (96%)	196 (97%)	6 (3%)	0	100	100
5	E	196/198 (99%)	188 (96%)	8 (4%)	0	100	100
6	F	406/414 (98%)	402 (99%)	4 (1%)	0	100	100
All	All	1882/1940 (97%)	1826 (97%)	56 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	376/376 (100%)	375 (100%)	1 (0%)	86	94
2	B	296/320 (92%)	292 (99%)	4 (1%)	59	76
3	C	200/205 (98%)	200 (100%)	0	100	100
4	D	171/176 (97%)	170 (99%)	1 (1%)	78	88
5	E	165/165 (100%)	164 (99%)	1 (1%)	78	88
6	F	337/343 (98%)	335 (99%)	2 (1%)	78	88
All	All	1545/1585 (98%)	1536 (99%)	9 (1%)	76	88

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

Mol	Chain	Res	Type
1	A	444	LYS
2	B	49	ASP
2	B	236	THR
2	B	382	MET
2	B	401	VAL
4	D	145	LEU
5	E	176	LEU
6	F	131	LYS
6	F	383	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	88	GLN
1	A	368	ASN
2	B	153	HIS
3	C	93	GLN
5	E	56	ASN
6	F	237	ASN

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)

Of 9 ligands modelled in this entry, 1 is monoatomic - leaving 8 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
9	LMT	D	301	-	36,36,36	0.58	0	47,47,47	0.79	0
7	FMN	C	301	-	33,33,33	0.63	0	48,50,50	0.71	1 (2%)
7	FMN	B	501	-	33,33,33	0.62	0	48,50,50	0.73	1 (2%)
11	FES	E	301	5,4	0,4,4	-	-	-		
12	FAD	F	502	-	56,58,58	0.82	1 (1%)	81,89,89	0.74	1 (1%)
11	FES	F	501	6	0,4,4	-	-	-		
9	LMT	B	503	-	36,36,36	0.60	0	47,47,47	0.72	0
8	RBF	B	502	-	29,29,29	1.01	1 (3%)	41,43,43	0.88	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
9	LMT	D	301	-	-	6/21/61/61	0/2/2/2

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
7	FMN	C	301	-	-	0/18/18/18	0/3/3/3
7	FMN	B	501	-	-	6/18/18/18	0/3/3/3
11	FES	E	301	5,4	-	-	0/1/1/1
12	FAD	F	502	-	-	8/34/50/50	0/6/6/6
11	FES	F	501	6	-	-	0/1/1/1
9	LMT	B	503	-	-	6/21/61/61	0/2/2/2
8	RBF	B	502	-	-	1/14/14/14	0/3/3/3

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
8	B	502	RBF	C5A-N5	-2.47	1.34	1.39
12	F	502	FAD	C5X-N5	-2.24	1.35	1.39

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
7	C	301	FMN	C4-N3-C2	-2.07	121.82	125.64
12	F	502	FAD	O2P-P-O1P	2.06	122.43	112.24
7	B	501	FMN	C4-N3-C2	-2.04	121.87	125.64

There are no chirality outliers.

All (27) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	B	501	FMN	C1'-C2'-C3'-C4'
7	B	501	FMN	C4'-C5'-O5'-P
9	B	503	LMT	C2'-C1'-O1'-C1
9	B	503	LMT	O5'-C1'-O1'-C1
12	F	502	FAD	N10-C1'-C2'-O2'
12	F	502	FAD	N10-C1'-C2'-C3'
12	F	502	FAD	O4'-C4'-C5'-O5'
9	D	301	LMT	C3'-C4'-O1B-C1B
7	B	501	FMN	O2'-C2'-C3'-O3'
7	B	501	FMN	O2'-C2'-C3'-C4'
9	D	301	LMT	C1-C2-C3-C4
9	D	301	LMT	O5B-C5B-C6B-O6B
9	B	503	LMT	O1'-C1-C2-C3
9	D	301	LMT	O5'-C5'-C6'-O6'
9	B	503	LMT	O5B-C5B-C6B-O6B

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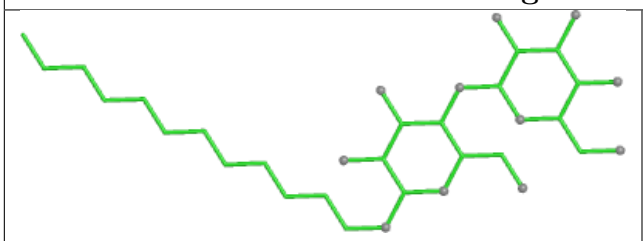
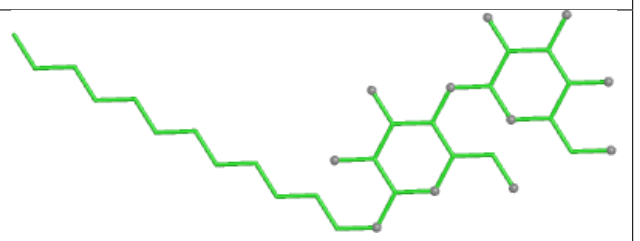
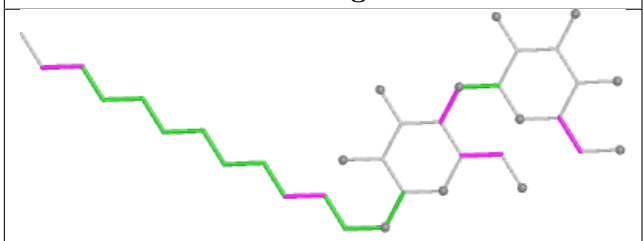
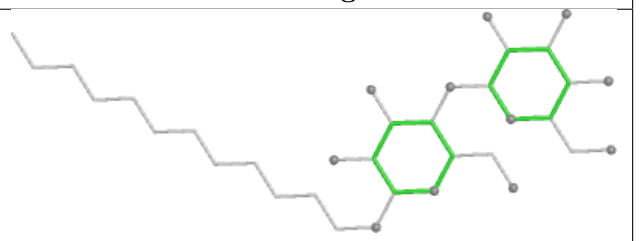
Mol	Chain	Res	Type	Atoms
8	B	502	RBF	C2'-C3'-C4'-O4'
12	F	502	FAD	C2'-C3'-C4'-O4'
9	B	503	LMT	C1-C2-C3-C4
12	F	502	FAD	C2'-C3'-C4'-C5'
12	F	502	FAD	O3'-C3'-C4'-O4'
12	F	502	FAD	C3'-C4'-C5'-O5'
12	F	502	FAD	P-O3P-PA-O2A
7	B	501	FMN	O4'-C4'-C5'-O5'
9	B	503	LMT	C6-C7-C8-C9
9	D	301	LMT	C5'-C4'-O1B-C1B
9	D	301	LMT	C9-C10-C11-C12
7	B	501	FMN	C1'-C2'-C3'-O3'

There are no ring outliers.

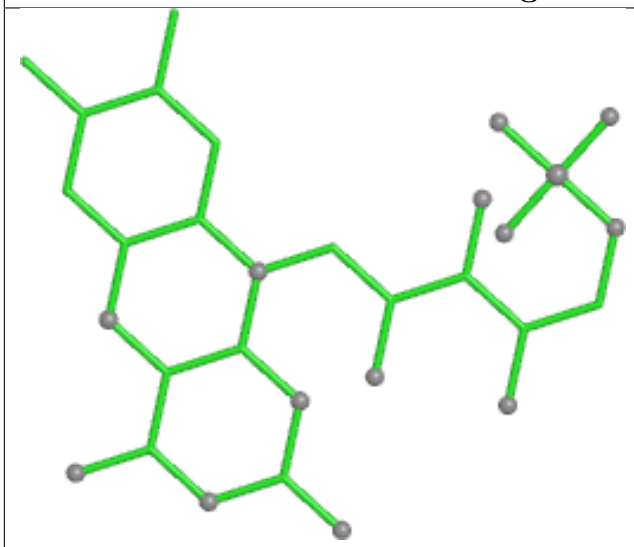
3 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
9	D	301	LMT	4	0
7	B	501	FMN	2	0
12	F	502	FAD	4	0

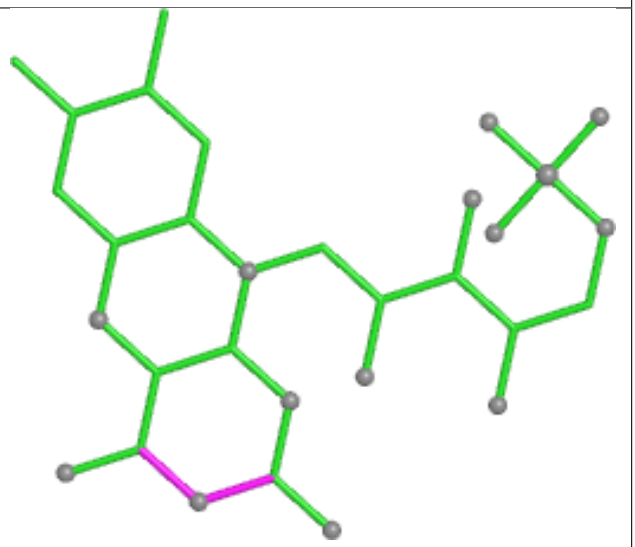
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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 LMT D 301	
 Bond lengths	 Bond angles
 Torsions	 Rings

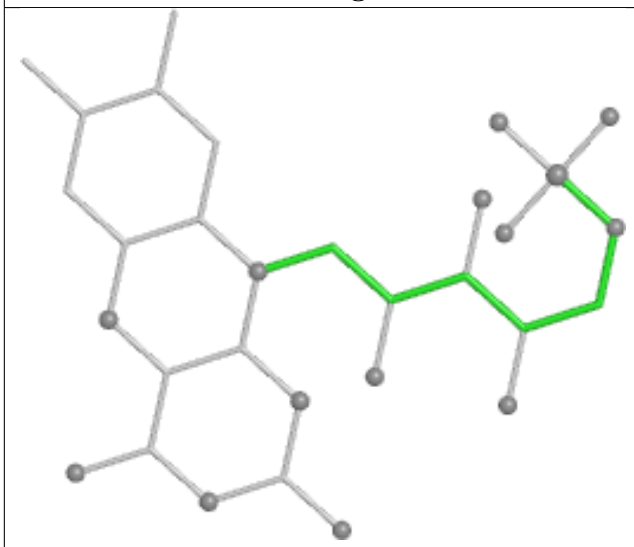
Ligand FMN C 301



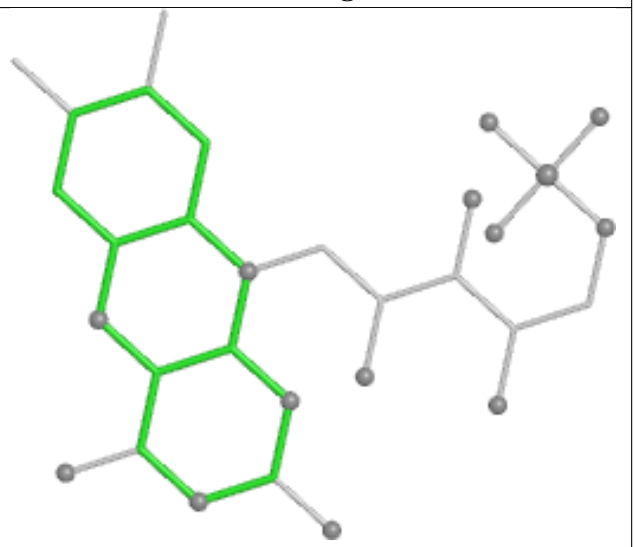
Bond lengths



Bond angles

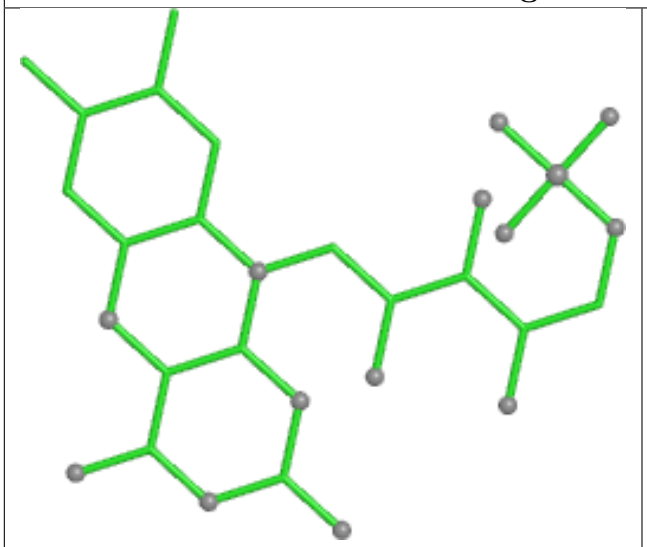


Torsions

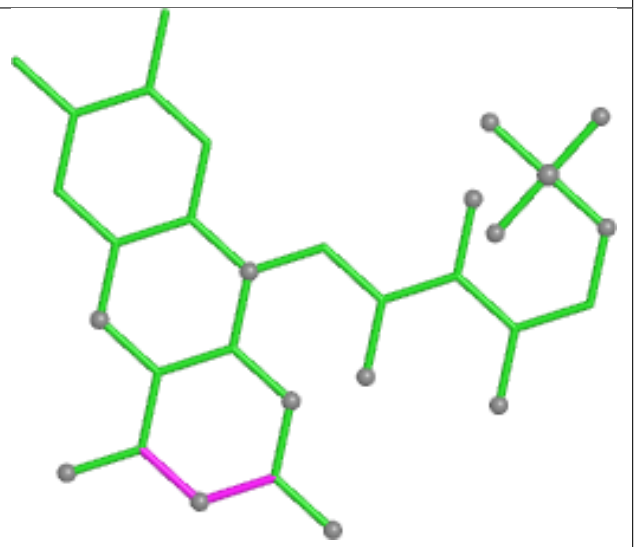


Rings

Ligand FMN B 501



Bond lengths



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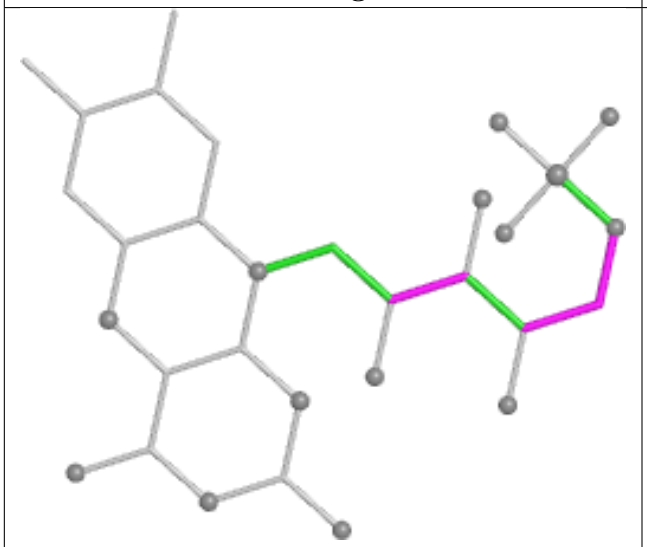
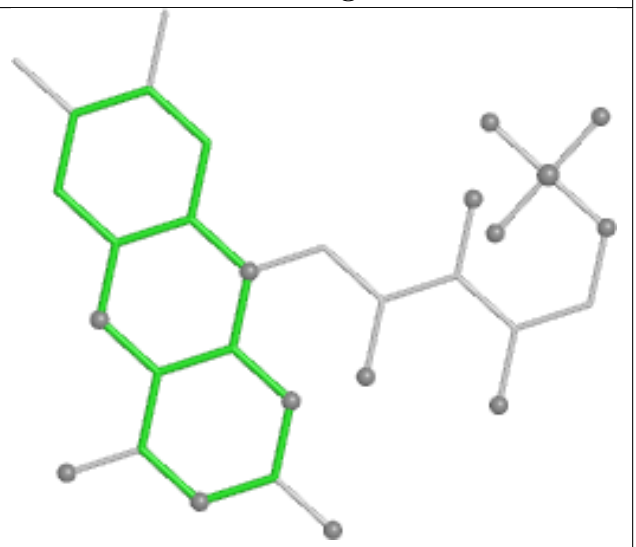
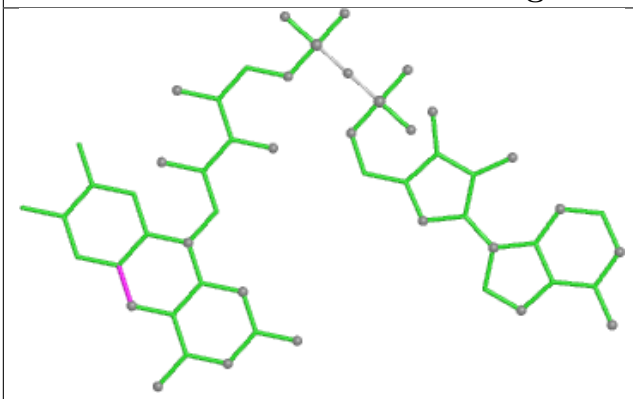


Torsions

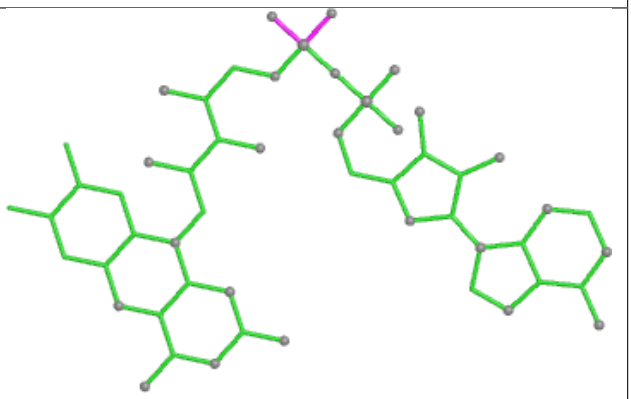


Rings

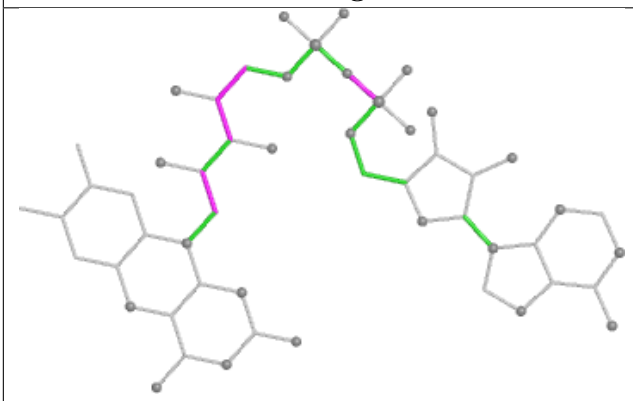
Ligand FAD F 502



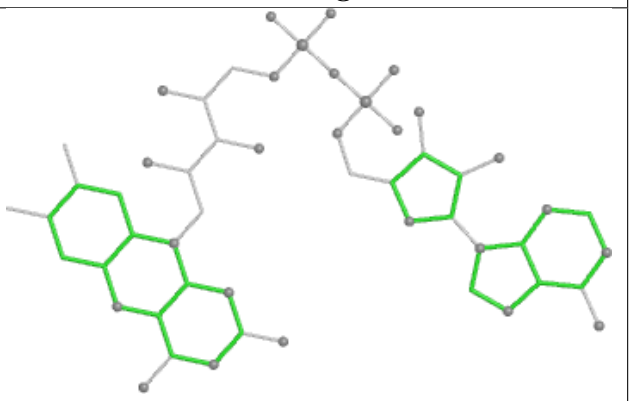
Bond lengths



Bond angles

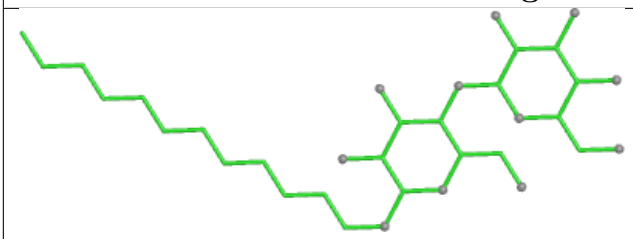


Torsions

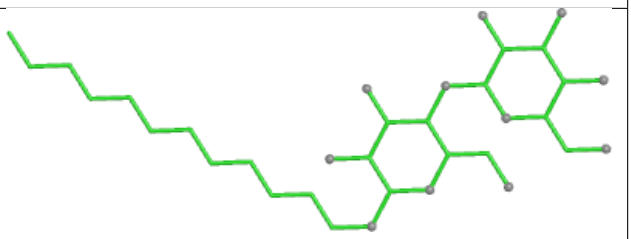


Rings

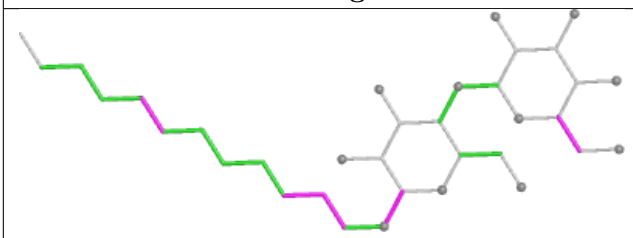
Ligand LMT B 503



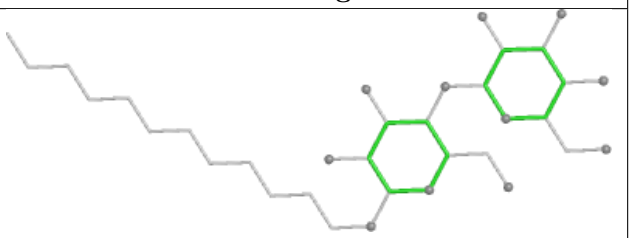
Bond lengths



Bond angles

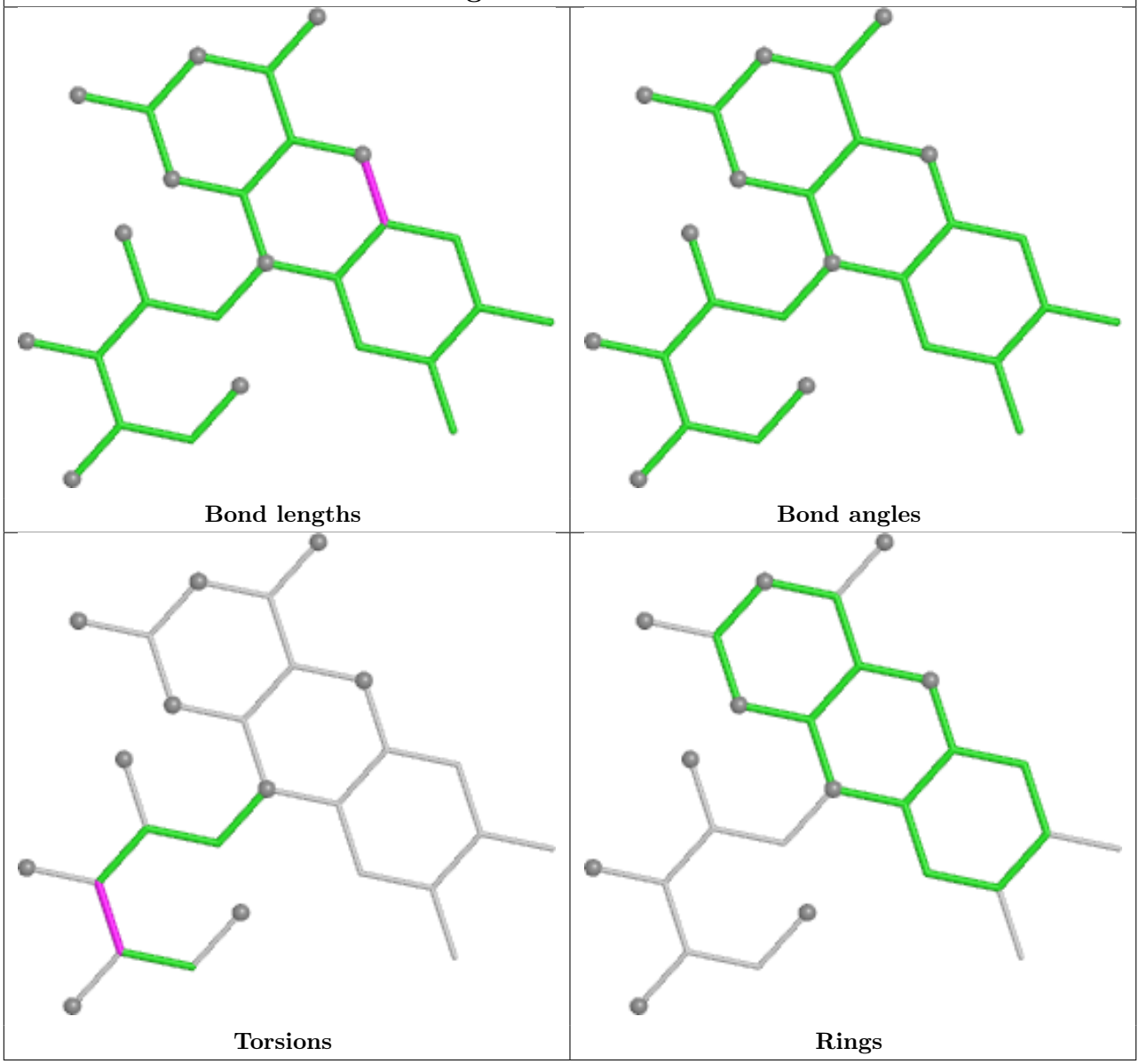


Torsions



Rings

Ligand RBF B 502



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.

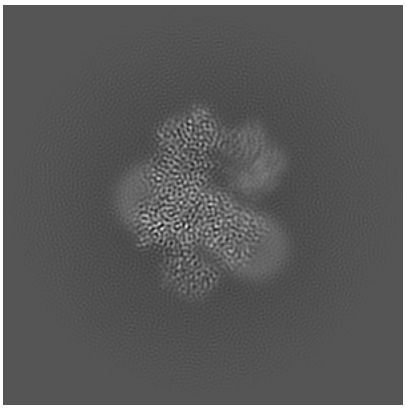
6 Map visualisation (i)

This section contains visualisations of the EMDB entry EMD-64063. These allow visual inspection of the internal detail of the map and identification of artifacts.

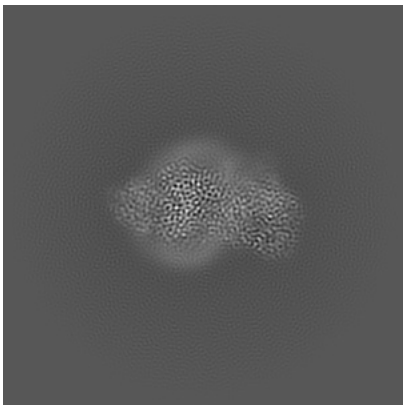
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections (i)

6.1.1 Primary map



X



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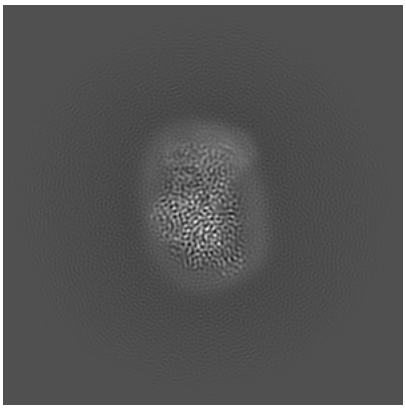
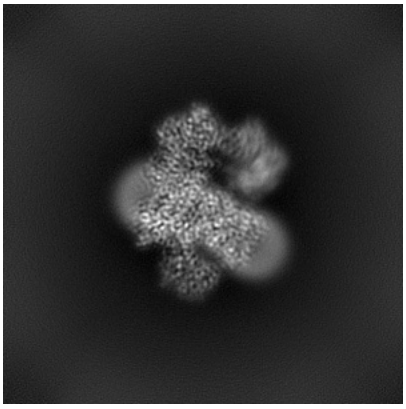


Z

6.1.2 Raw map



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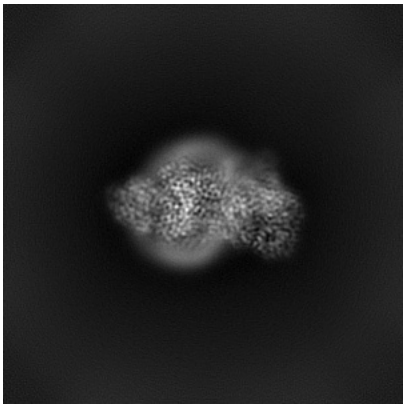

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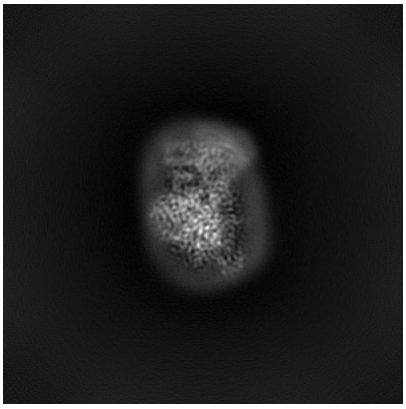
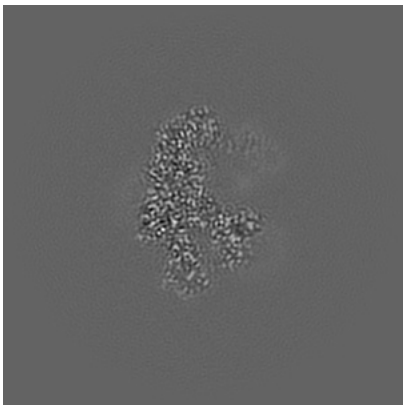


Z

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

6.2 Central slices (i)

6.2.1 Primary map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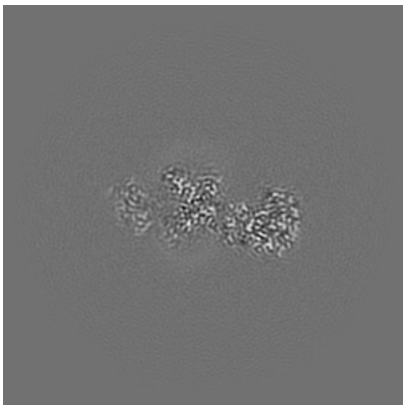
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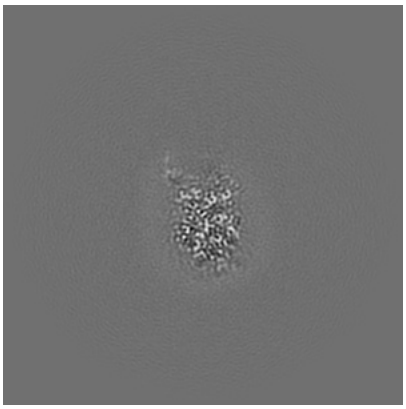
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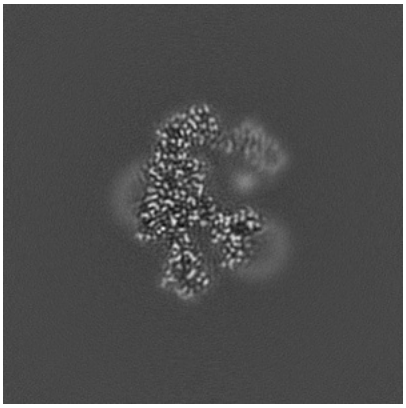
6.2.2 Raw map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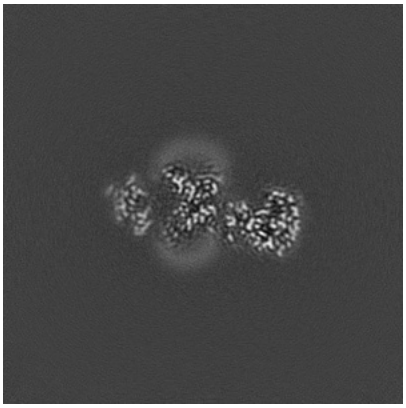
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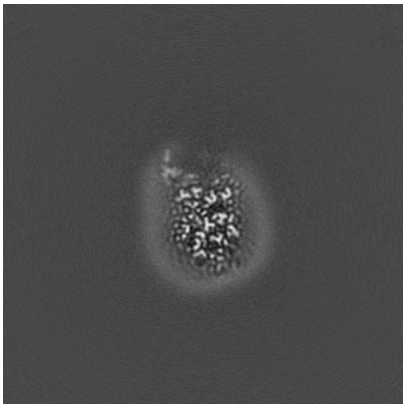
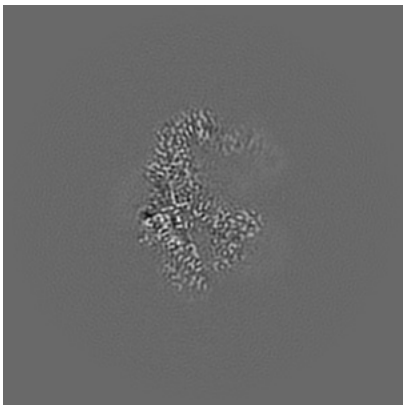
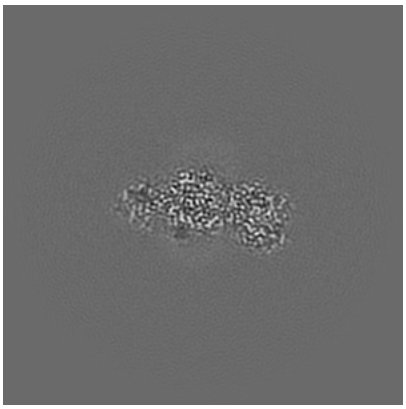

Z Index: 160

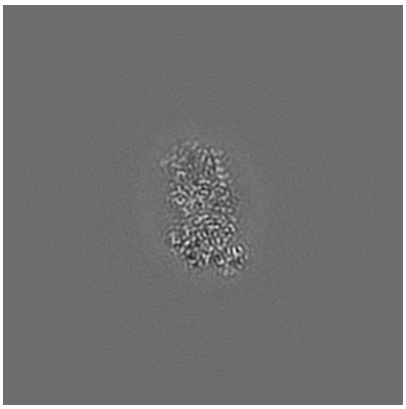
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices (i)

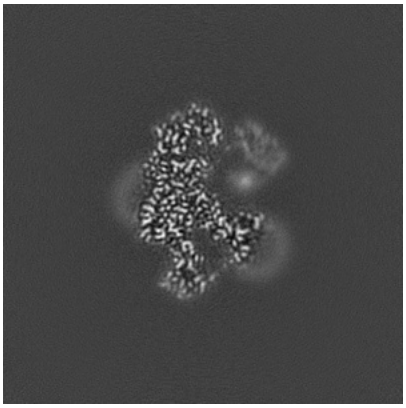
6.3.1 Primary map

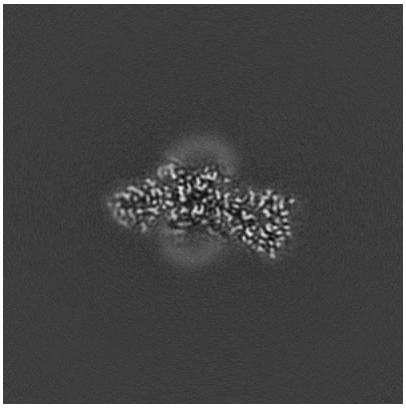

X Index: 163

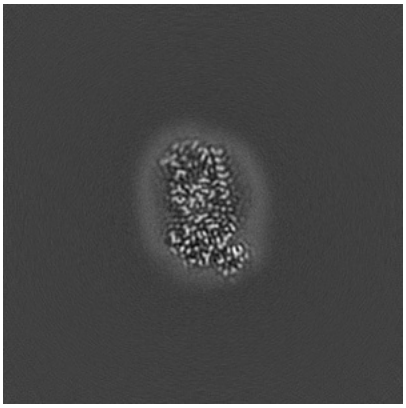

Y Index: 139


Z Index: 144

6.3.2 Raw map


X Index: 157

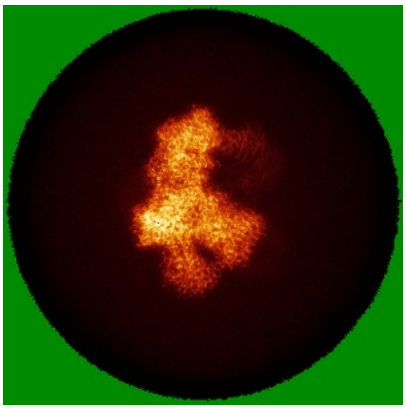

Y Index: 146

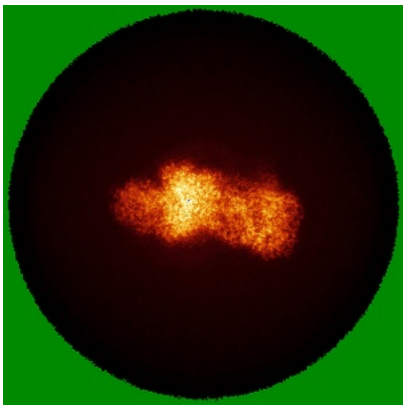

Z Index: 144

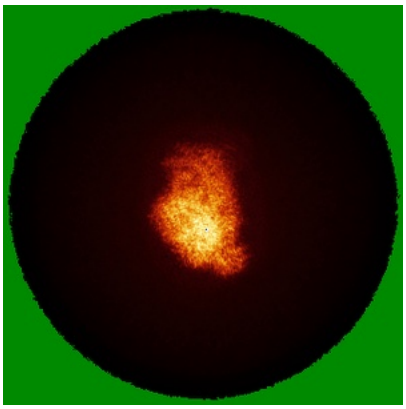
The images above show the largest variance slices of the map in three orthogonal directions.

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

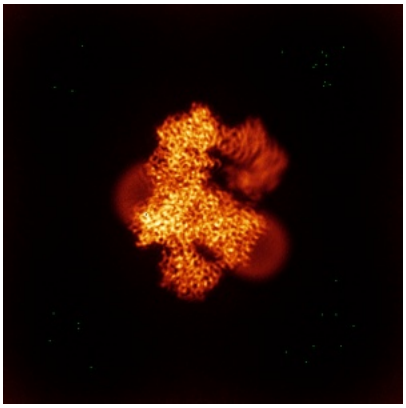
6.4.1 Primary map

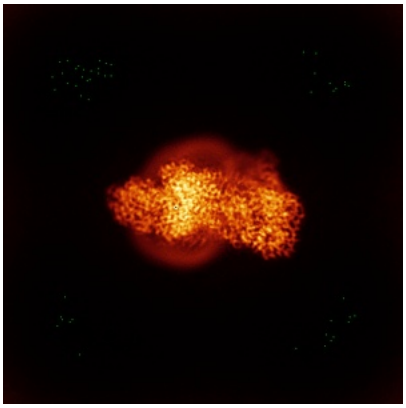

X

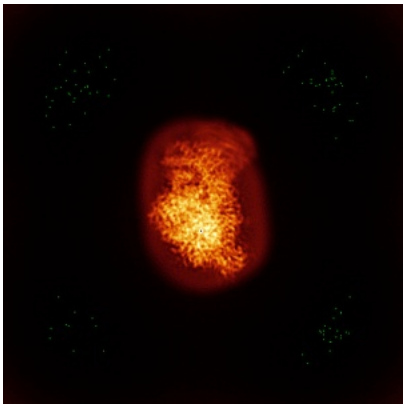

Y


Z

6.4.2 Raw map


X


Y


Z

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

6.5 Orthogonal surface views (i)

6.5.1 Primary map

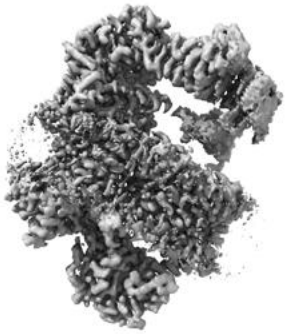
X

Y

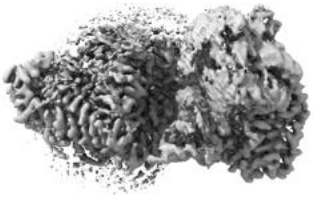
Z

The images above show the 3D surface view of the map at the recommended contour level 0.195. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.5.2 Raw map



X



Y

Z

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

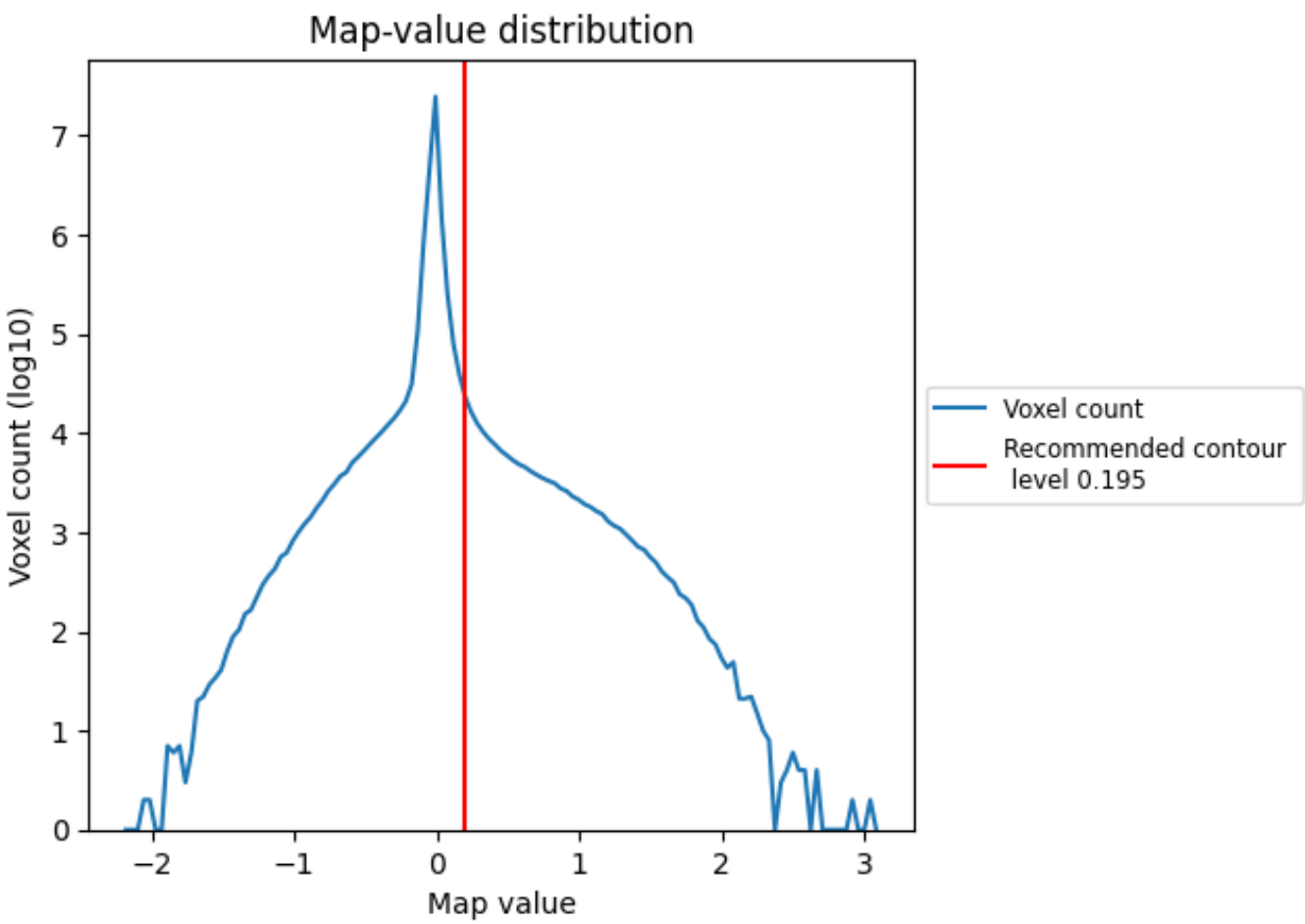
6.6 Mask visualisation (i)

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

7 Map analysis (i)

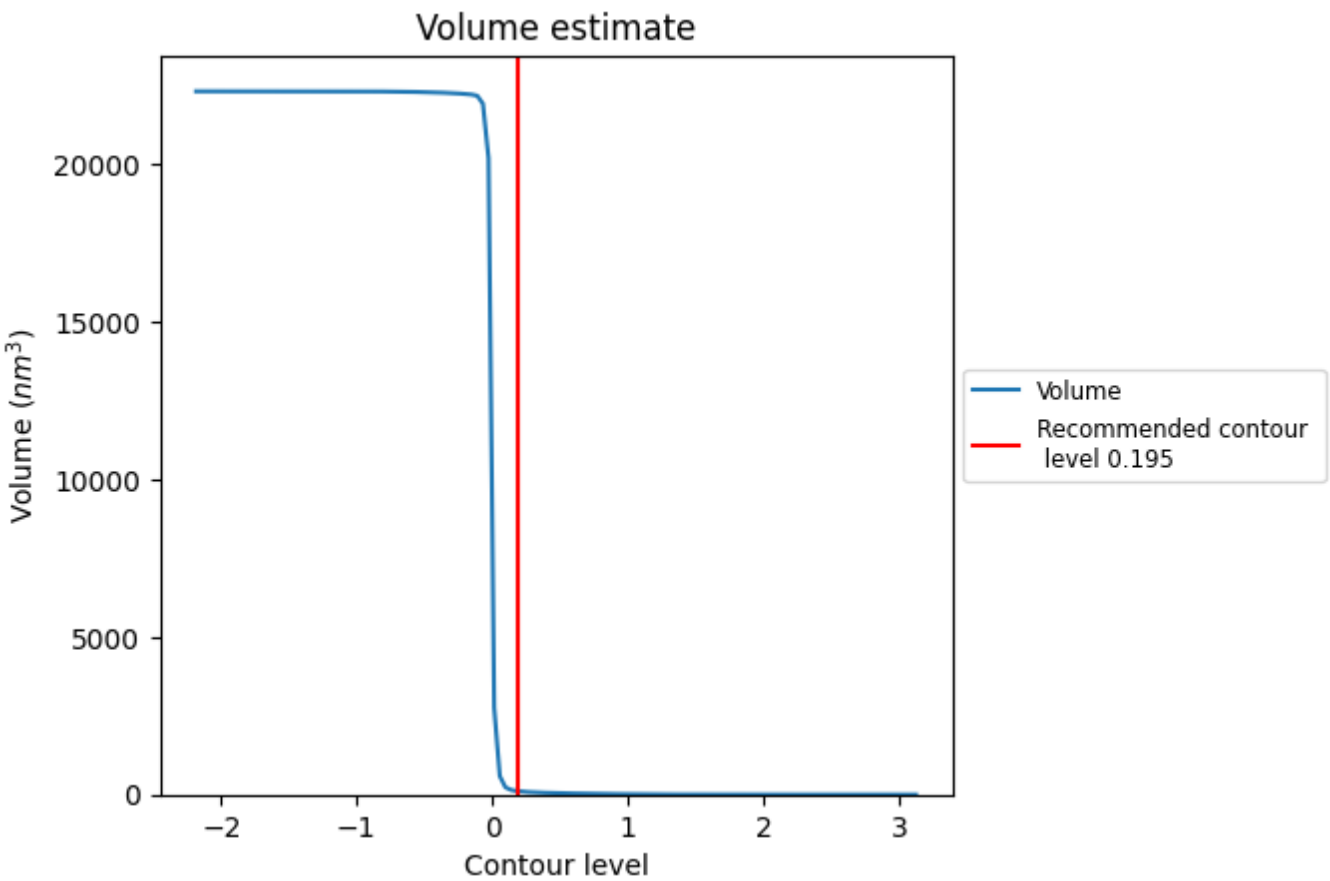
This section contains the results of statistical analysis of the map.

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.

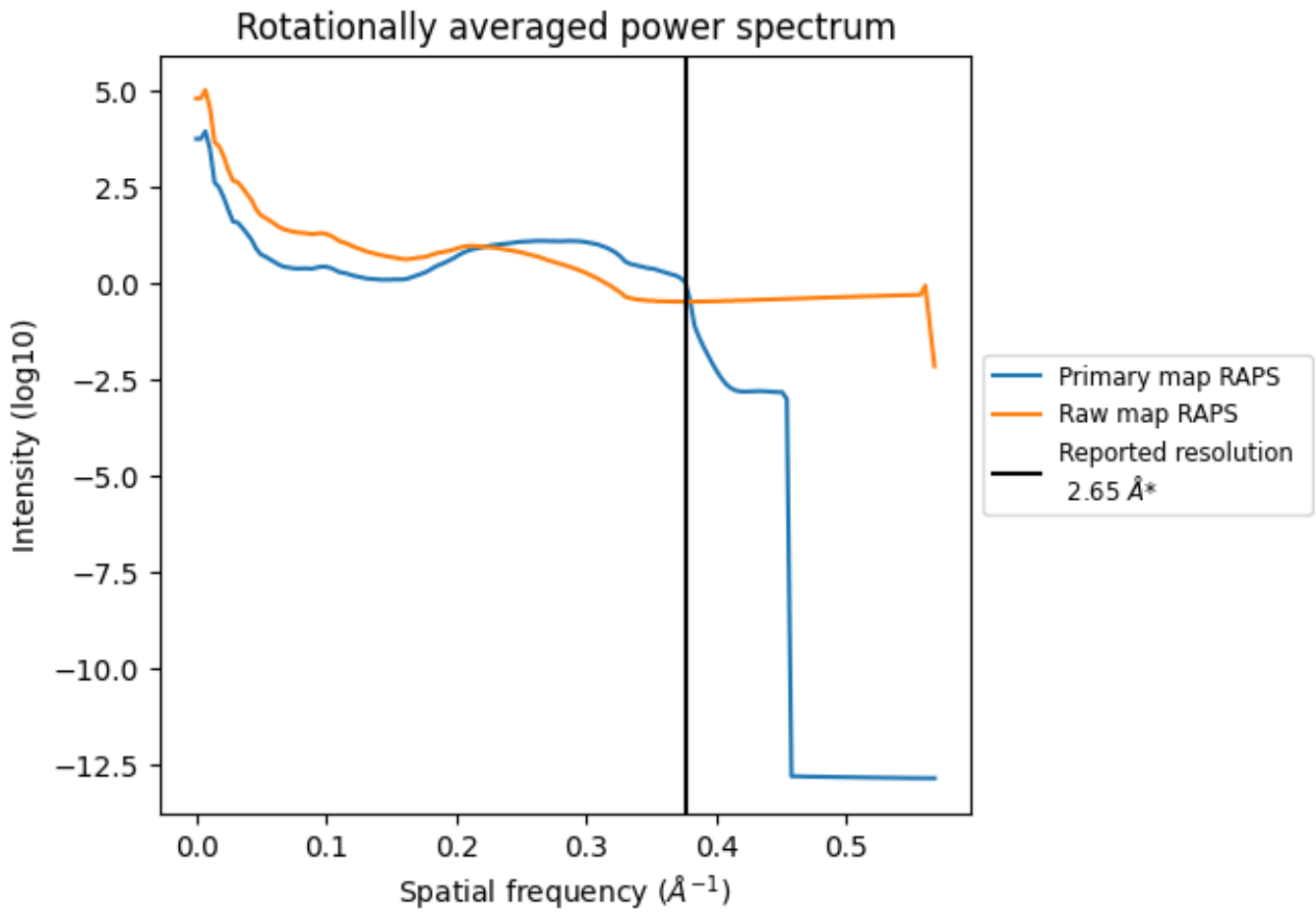
7.2 Volume estimate (i)



The volume at the recommended contour level is 106 nm^3; this corresponds to an approximate mass of 96 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)

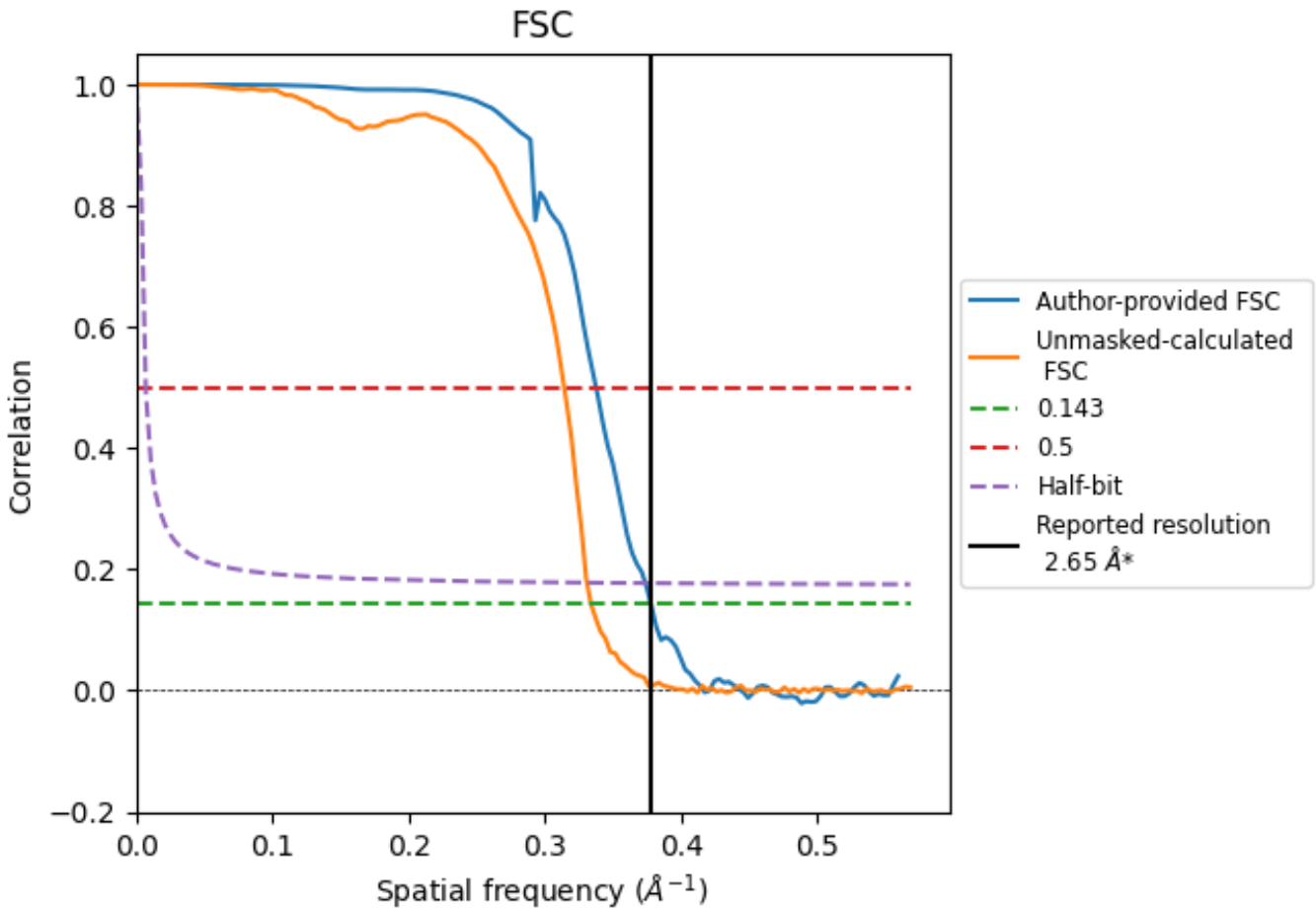


*Reported resolution corresponds to spatial frequency of 0.377 Å^{-1}

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 0.377 Å^{-1}

8.2 Resolution estimates (i)

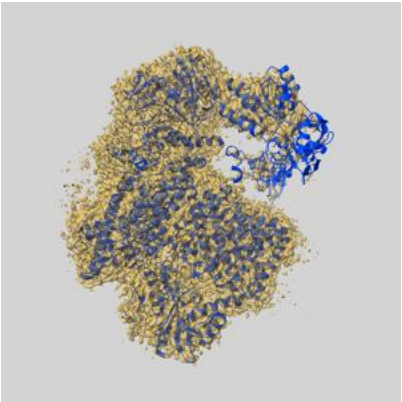
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.65	-	-
Author-provided FSC curve	2.65	2.96	2.68
Unmasked-calculated*	3.00	3.18	3.02

*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 3.00 differs from the reported value 2.65 by more than 10 %

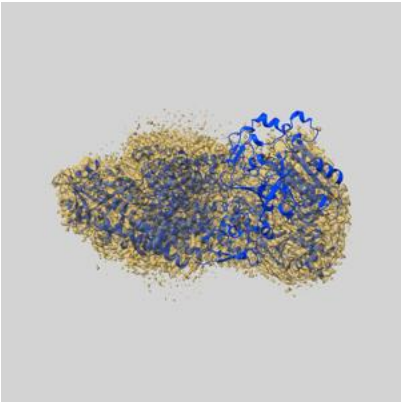
9 Map-model fit (i)

This section contains information regarding the fit between EMDB map EMD-64063 and PDB model 9UD6. Per-residue inclusion information can be found in section 3 on page 9.

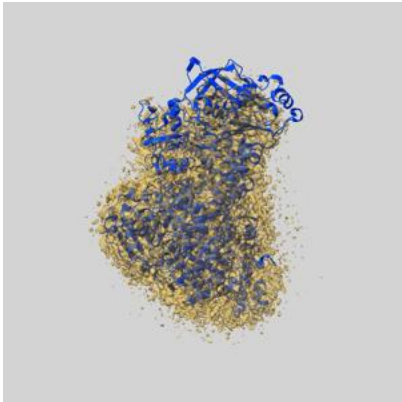
9.1 Map-model overlay (i)



X



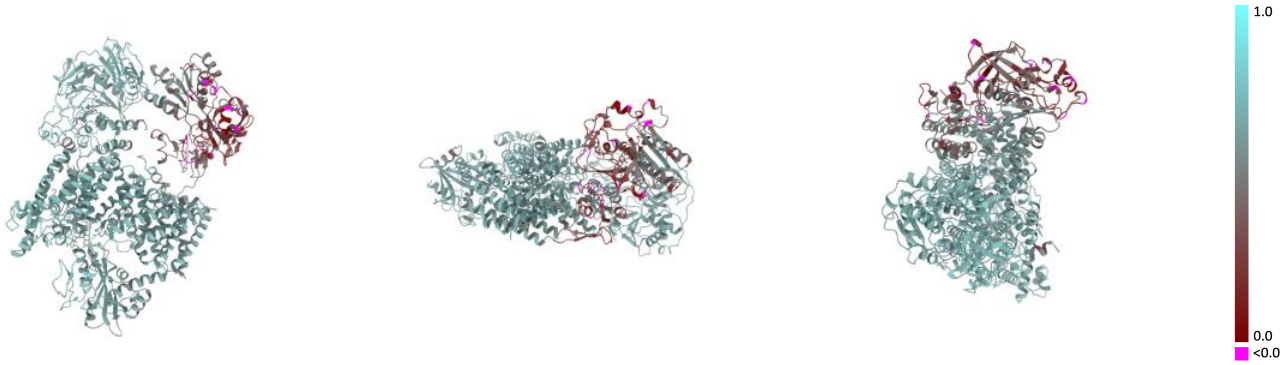
Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.195 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.

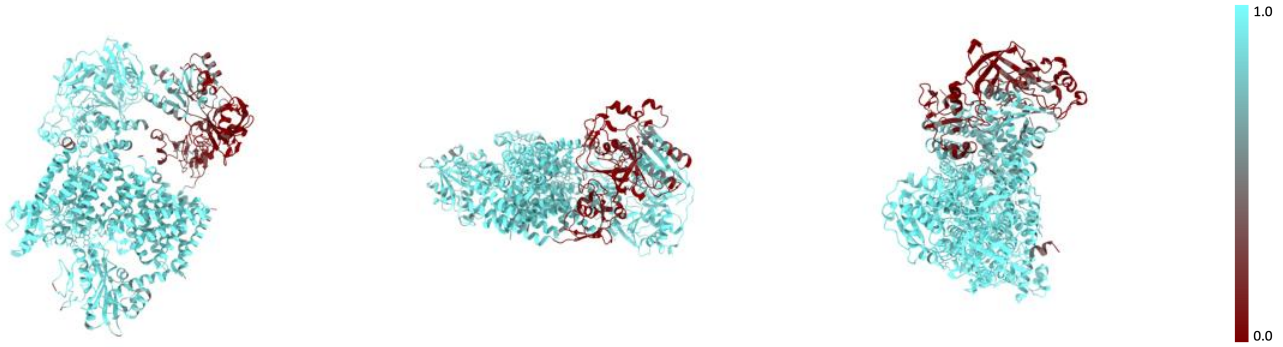
9.2 Q-score mapped to coordinate model (i)



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.

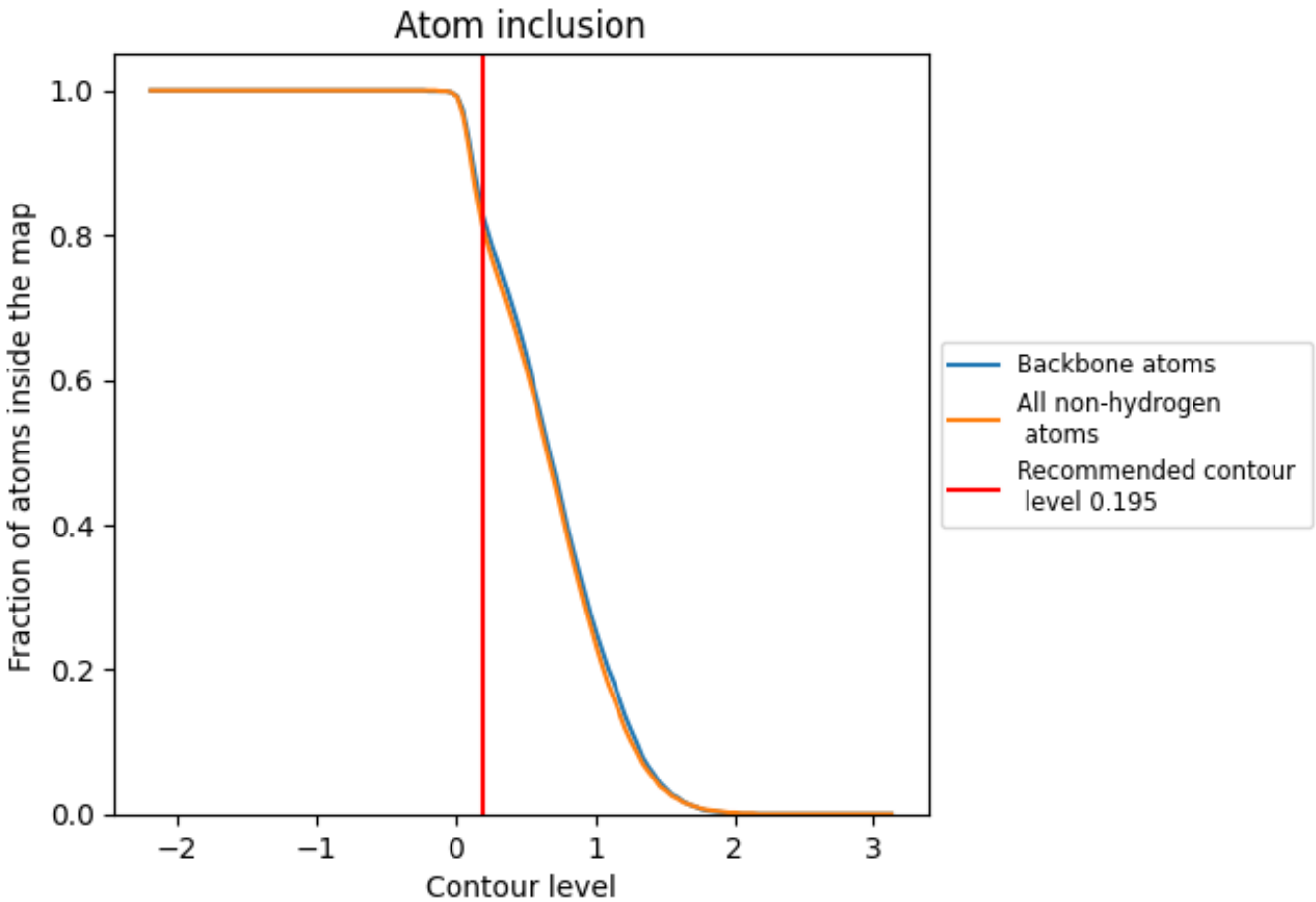
9.3 Atom inclusion mapped to coordinate model (i)



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 (0.195).

9.4 Atom inclusion (i)



At the recommended contour level, 83% of all backbone atoms, 81% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary (i)

The table lists the average atom inclusion at the recommended contour level (0.195) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.8120	0.5700
A	0.9620	0.6400
B	0.9760	0.6500
C	0.8970	0.6050
D	0.9560	0.6280
E	0.9520	0.6260
F	0.3120	0.3440