



Full wwPDB EM Validation Report ⓘ

Nov 3, 2024 – 07:16 AM EST

PDB ID : 6WXV
EMDB ID : EMD-21964
Title : CryoEM structure of mouse DUOX1-DUOXA1 complex in the presence of NADPH
Authors : Sun, J.
Deposited on : 2020-05-12
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 : 2022.3.0, CSD as543be (2022)
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

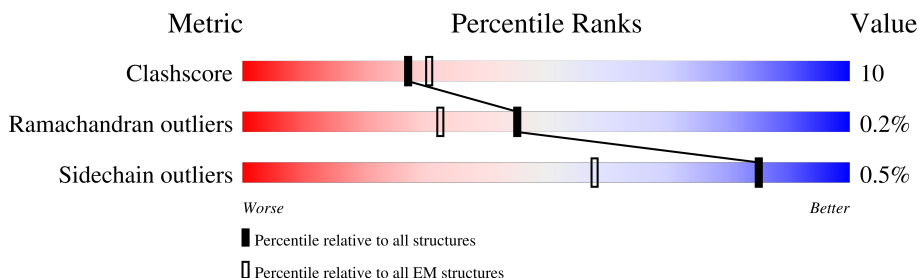
1 Overall quality at a glance

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	1536	
2	B	341	
3	C	5	

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
6	HEB	A	1605	X	-	-	-
6	HEB	A	1606	X	-	-	-

2 Entry composition [i](#)

There are 9 unique types of molecules in this entry. The entry contains 10605 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 Dual oxidase 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	1091	Total	C	N	O	S	0	0
			8524	5538	1485	1470	31		

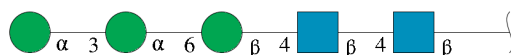
There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	16	GLY	-	expression tag	UNP A2AQ92
A	17	PRO	-	expression tag	UNP A2AQ92
A	18	SER	-	expression tag	UNP A2AQ92
A	19	ARG	-	expression tag	UNP A2AQ92

- Molecule 2 is a protein called Dual oxidase maturation factor 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	228	Total	C	N	O	S	0	0
			1742	1150	289	293	10		

- Molecule 3 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-alpha-D-mannopyranose-(1-6)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



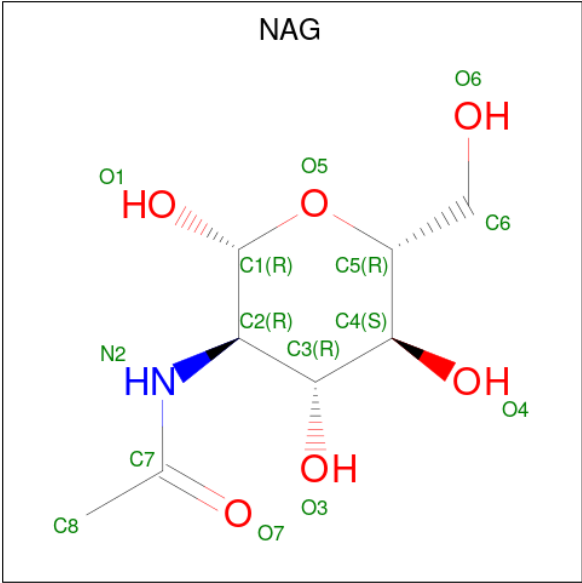
Mol	Chain	Residues	Atoms				AltConf	Trace
3	C	5	Total	C	N	O	0	0
			61	34	2	25		

- Molecule 4 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula: C₂₇H₃₃N₉O₁₅P₂).



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

- Molecule 5 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: C₈H₁₅NO₆).



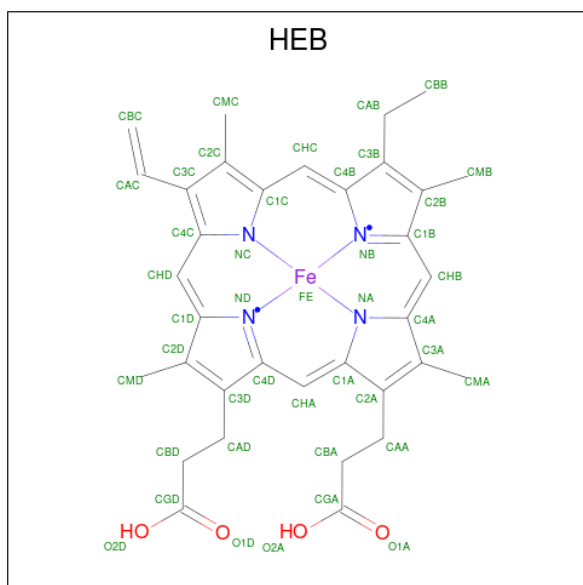
Mol	Chain	Residues	Atoms				AltConf
5	A	1	Total	C	N	O	0
			14	8	1	5	
5	A	1	Total	C	N	O	0
			14	8	1	5	

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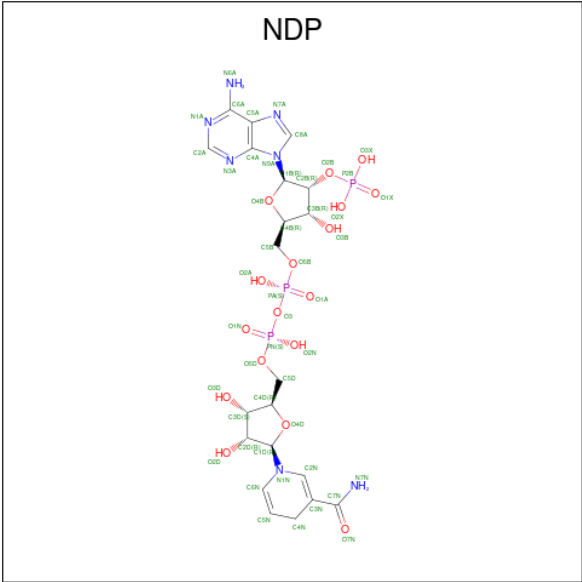
Mol	Chain	Residues	Atoms				AltConf
5	A	1	Total	C	N	O	0
			14	8	1	5	
5	B	1	Total	C	N	O	0
			14	8	1	5	

- Molecule 6 is HEME B/C (three-letter code: HEB) (formula: $C_{34}H_{34}FeN_4O_4$).



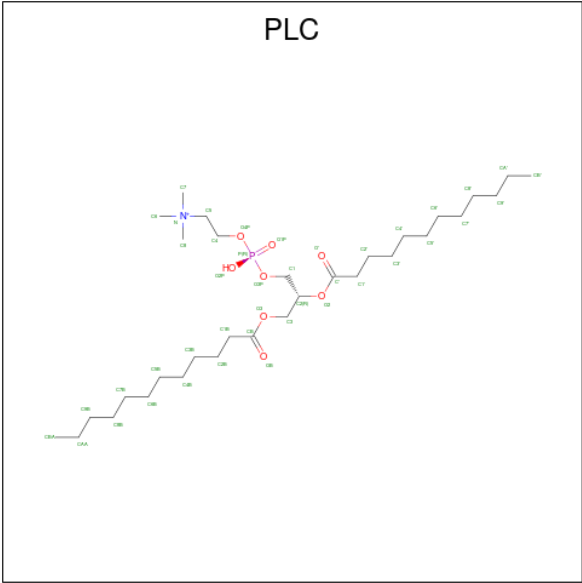
Mol	Chain	Residues	Atoms					AltConf
6	A	1	Total	C	Fe	N	O	0
			43	34	1	4	4	
6	A	1	Total	C	Fe	N	O	0
			43	34	1	4	4	

- Molecule 7 is NADPH DIHYDRO-NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (three-letter code: NDP) (formula: $C_{21}H_{30}N_7O_{17}P_3$) (labeled as "Ligand of Interest" by depositor).

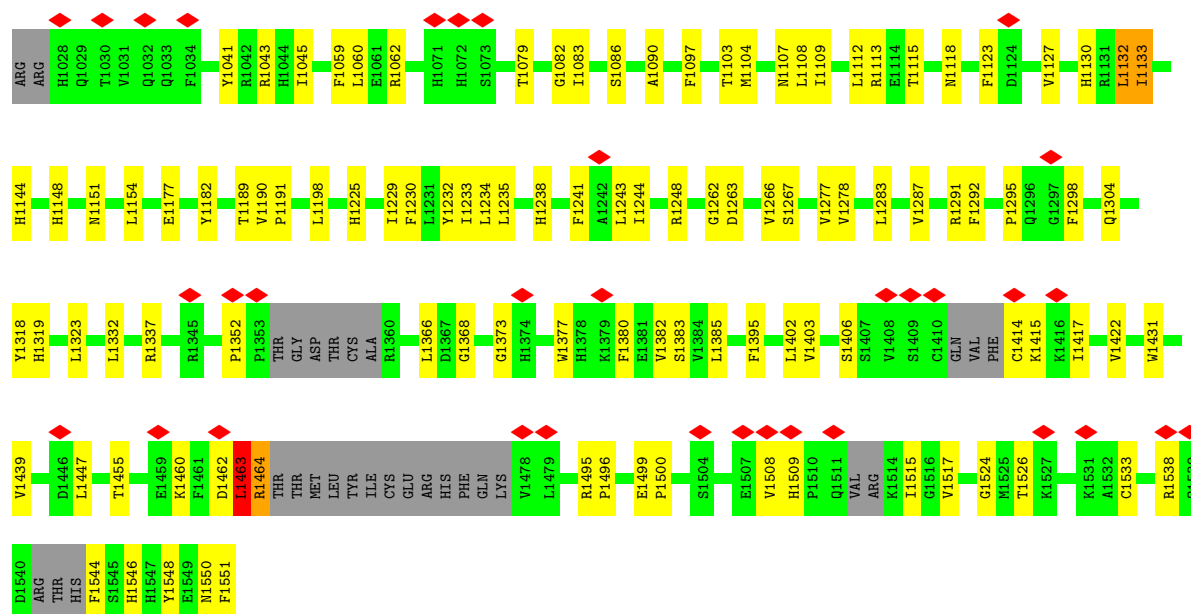


Mol	Chain	Residues	Atoms					AltConf
7	A	1	Total	C	N	O	P	0
			40	15	6	16	3	

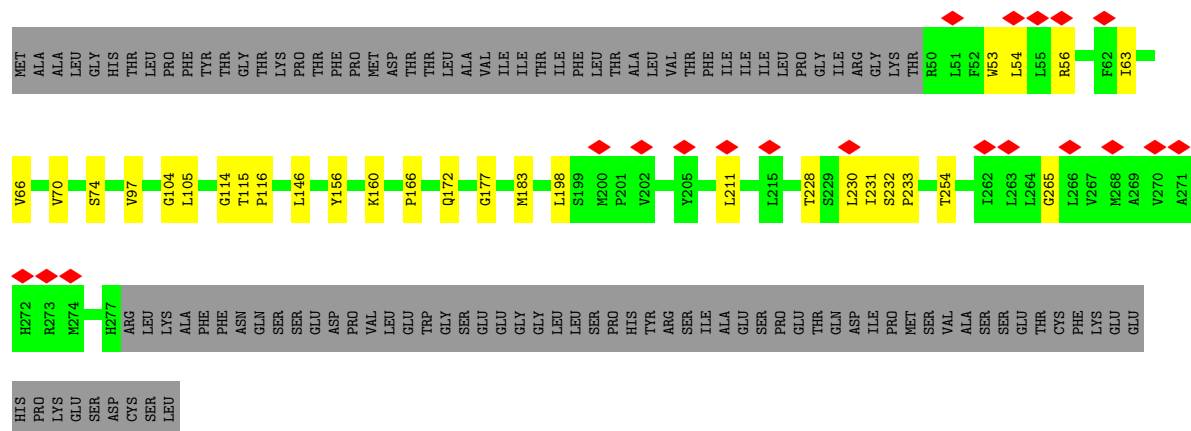
- Molecule 8 is DIUNDECYL PHOSPHATIDYL CHOLINE (three-letter code: PLC) (formula: $C_{32}H_{65}NO_8P$) (labeled as "Ligand of Interest" by depositor).



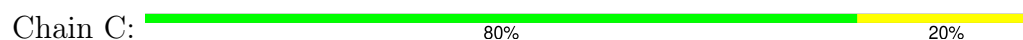
Mol	Chain	Residues	Atoms		AltConf
9	A	1	Total 1	Ca 1	0



• Molecule 2: Dual oxidase maturation factor 1



• Molecule 3: alpha-D-mannopyranose-(1-3)-alpha-D-mannopyranose-(1-6)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	667038	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 ($e^-/\text{\AA}^2$)	65.4	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	3.943	Depositor
Minimum map value	-1.874	Depositor
Average map value	0.011	Depositor
Map value standard deviation	0.095	Depositor
Recommended contour level	0.55	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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: FAD, NDP, MAN, BMA, HEB, CA, PLC, NAG

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	0.35	1/8784 (0.0%)	0.48	5/11967 (0.0%)
2	B	0.26	0/1793	0.45	0/2455
All	All	0.34	1/10577 (0.0%)	0.48	5/14422 (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	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	1133	ILE	C-O	-5.29	1.13	1.23

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	214	LEU	CA-CB-CG	6.46	130.16	115.30
1	A	1133	ILE	CG1-CB-CG2	-6.09	98.00	111.40
1	A	1132	LEU	CB-CG-CD1	-5.96	100.88	111.00
1	A	213	LEU	CB-CG-CD2	-5.94	100.91	111.00
1	A	1133	ILE	CB-CA-C	-5.11	101.37	111.60

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	118	CYS	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	8524	0	8214	168	0
2	B	1742	0	1742	19	0
3	C	61	0	52	0	0
4	A	53	0	31	3	0
5	A	42	0	39	1	0
5	B	14	0	13	0	0
6	A	86	0	64	7	0
7	A	40	0	19	8	0
8	A	42	0	64	11	0
9	A	1	0	0	1	0
All	All	10605	0	10238	199	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 10.

All (199) 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)
7:A:1607:NDP:O3D	8:A:1608:PLC:H81	1.59	1.00
7:A:1607:NDP:H52N	8:A:1608:PLC:H71	1.55	0.88
1:A:213:LEU:H	1:A:213:LEU:HD23	1.40	0.85
4:A:1601:FAD:H9	4:A:1601:FAD:O3'	1.77	0.84
1:A:1189:THR:HG22	1:A:1191:PRO:HD2	1.63	0.79
1:A:1104:MET:CE	1:A:1127:VAL:HG22	2.14	0.78
1:A:1104:MET:HE1	1:A:1127:VAL:HG22	1.66	0.78
1:A:213:LEU:HD23	1:A:213:LEU:N	1.99	0.78
4:A:1601:FAD:O4'	4:A:1601:FAD:O2'	2.00	0.77
1:A:1403:VAL:HG21	1:A:1439:VAL:HG13	1.67	0.76
2:B:172:GLN:HE22	2:B:230:LEU:HB3	1.52	0.74
1:A:213:LEU:HD12	2:B:166:PRO:O	1.87	0.74
1:A:218:ALA:O	1:A:370:ARG:NH2	2.21	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:97:VAL:HA	2:B:114:GLY:HA2	1.71	0.70
1:A:227:GLY:O	1:A:370:ARG:NH1	2.24	0.70
1:A:66:PRO:HD3	1:A:148:ARG:HH12	1.56	0.69
1:A:289:ASN:ND2	1:A:532:ILE:O	2.25	0.69
1:A:83:MET:O	1:A:405:GLN:NE2	2.25	0.68
1:A:1464:ARG:O	1:A:1464:ARG:HG3	1.94	0.67
1:A:307:TYR:HA	1:A:509:ARG:HH12	1.60	0.66
1:A:1107:ASN:ND2	1:A:1267:SER:OG	2.28	0.66
6:A:1605:HEB:HHA	6:A:1605:HEB:HBA2	1.76	0.66
1:A:82:VAL:HG23	1:A:83:MET:HG3	1.76	0.66
1:A:341:ARG:NH1	1:A:565:CYS:SG	2.69	0.65
7:A:1607:NDP:O3D	8:A:1608:PLC:C8	2.41	0.65
1:A:176:SER:OG	9:A:1609:CA:CA	1.74	0.65
1:A:36:TYR:OH	2:B:146:LEU:O	2.13	0.64
1:A:1082:GLY:O	1:A:1151:ASN:ND2	2.30	0.63
1:A:315:ASP:O	1:A:507:ARG:NH2	2.31	0.63
1:A:571:LEU:HD21	1:A:576:LEU:HD11	1.81	0.63
1:A:1045:ILE:HG12	8:A:1608:PLC:H4'2	1.79	0.62
1:A:541:LEU:O	1:A:545:THR:OG1	2.14	0.62
1:A:306:GLU:O	1:A:509:ARG:NH2	2.27	0.61
1:A:1380:PHE:O	1:A:1414:CYS:N	2.34	0.61
1:A:1517:VAL:HB	1:A:1546:HIS:HA	1.82	0.61
1:A:1234:LEU:O	1:A:1238:HIS:HB2	2.00	0.61
1:A:206:PRO:HG2	1:A:232:TYR:HB3	1.81	0.60
1:A:472:TYR:OH	1:A:480:GLU:OE1	2.15	0.60
1:A:1277:VAL:HA	1:A:1292:PHE:HB3	1.83	0.60
1:A:188:THR:OG1	1:A:214:LEU:HD11	2.02	0.60
1:A:1133:ILE:HG22	1:A:1133:ILE:O	2.01	0.60
1:A:1283:LEU:HB2	1:A:1287:VAL:HG23	1.84	0.59
1:A:1422:VAL:HG11	1:A:1496:PRO:HG3	1.84	0.59
2:B:53:TRP:HE3	2:B:54:LEU:HD12	1.67	0.59
1:A:76:ARG:NH1	1:A:418:ASP:OD2	2.35	0.59
1:A:79:SER:O	1:A:405:GLN:NE2	2.37	0.58
1:A:208:ASP:C	1:A:211:SER:OG	2.43	0.57
1:A:585:ARG:NH1	1:A:1177:GLU:OE2	2.37	0.57
1:A:190:ARG:NH2	1:A:273:GLU:OE1	2.32	0.57
1:A:1043:ARG:NH2	1:A:1118:ASN:O	2.38	0.57
1:A:213:LEU:CD1	2:B:166:PRO:O	2.53	0.57
1:A:119:PRO:HD3	1:A:153:ARG:HH12	1.70	0.57
1:A:418:ASP:HB3	1:A:421:ALA:HB3	1.86	0.57
5:A:1603:NAG:H3	5:A:1603:NAG:H83	1.86	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:212:SER:HA	1:A:217:MET:HE1	1.87	0.56
1:A:333:MET:HA	1:A:394:GLU:HB2	1.88	0.56
1:A:1406:SER:HB2	1:A:1447:LEU:HD13	1.87	0.56
1:A:504:GLN:HE21	1:A:508:LEU:HG	1.71	0.55
1:A:170:THR:HG23	1:A:181:SER:HB2	1.87	0.55
1:A:1229:ILE:O	1:A:1233:ILE:HG13	2.07	0.55
1:A:1298:PHE:HE1	1:A:1368:GLY:HA2	1.72	0.55
1:A:1241:PHE:HD2	1:A:1243:LEU:HG	1.72	0.54
1:A:1132:LEU:O	1:A:1132:LEU:HD12	2.07	0.54
1:A:1086:SER:HG	1:A:1148:HIS:CD2	2.22	0.54
1:A:1079:THR:HG22	1:A:1154:LEU:HD12	1.88	0.54
1:A:513:ARG:O	1:A:521:ASN:ND2	2.36	0.54
1:A:1262:GLY:O	1:A:1266:VAL:HG23	2.09	0.53
1:A:339:TYR:CE1	1:A:571:LEU:HD12	2.43	0.53
6:A:1605:HEB:HMB1	6:A:1605:HEB:HBB3	1.91	0.53
2:B:74:SER:HA	2:B:104:GLY:HA3	1.89	0.53
1:A:208:ASP:C	1:A:211:SER:HG	2.11	0.53
1:A:1385:LEU:HB3	1:A:1395:PHE:HE2	1.75	0.52
1:A:382:ASP:OD1	1:A:545:THR:HB	2.10	0.52
1:A:1104:MET:HE2	1:A:1127:VAL:HG22	1.90	0.51
1:A:1533:CYS:SG	1:A:1544:PHE:HB3	2.50	0.51
1:A:1238:HIS:HE1	6:A:1606:HEB:NA	2.08	0.51
1:A:117:GLY:HA3	1:A:150:ARG:HA	1.93	0.51
1:A:307:TYR:HD1	1:A:509:ARG:HH11	1.59	0.51
1:A:250:LEU:HD12	1:A:381:VAL:HG21	1.93	0.51
6:A:1606:HEB:HMB1	6:A:1606:HEB:HBB2	1.91	0.51
1:A:208:ASP:O	1:A:211:SER:OG	2.24	0.50
1:A:427:GLY:HA2	1:A:432:LEU:HD23	1.94	0.50
1:A:91:SER:OG	1:A:396:GLU:OE2	2.27	0.50
1:A:1103:THR:HG21	1:A:1130:HIS:HB2	1.93	0.50
1:A:337:GLY:H	1:A:571:LEU:HD22	1.76	0.50
1:A:51:LEU:HD11	1:A:113:VAL:HG11	1.94	0.50
1:A:1366:LEU:HD23	1:A:1368:GLY:H	1.77	0.50
2:B:70:VAL:HG11	2:B:254:THR:OG1	2.11	0.49
1:A:149:SER:HB3	1:A:426:ARG:HH12	1.77	0.49
1:A:1190:VAL:HB	1:A:1191:PRO:HD3	1.94	0.49
1:A:612:SER:O	1:A:616:VAL:HG23	2.13	0.49
1:A:1112:LEU:O	1:A:1115:THR:OG1	2.30	0.49
1:A:1508:VAL:HG23	1:A:1509:HIS:ND1	2.28	0.49
2:B:211:LEU:HD22	2:B:265:GLY:HA3	1.95	0.49
1:A:1097:PHE:O	1:A:1232:TYR:OH	2.23	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:504:GLN:NE2	1:A:508:LEU:HG	2.28	0.49
2:B:105:LEU:HD22	2:B:177:GLY:HA2	1.95	0.48
1:A:468:THR:HA	1:A:471:LEU:HD12	1.95	0.48
7:A:1607:NDP:H52N	8:A:1608:PLC:C7	2.36	0.48
1:A:362:ARG:NE	1:A:364:CYS:SG	2.84	0.48
2:B:53:TRP:CE3	2:B:54:LEU:HD12	2.48	0.48
1:A:341:ARG:NE	1:A:386:LEU:O	2.47	0.48
1:A:433:PRO:HB3	1:A:437:LYS:HE3	1.95	0.48
1:A:129:HIS:ND1	1:A:140:GLY:O	2.40	0.47
1:A:54:LEU:HD23	1:A:442:LEU:HD21	1.95	0.47
1:A:361:LEU:H	1:A:361:LEU:HD23	1.78	0.47
1:A:503:ASP:OD1	1:A:504:GLN:N	2.48	0.47
1:A:1198:LEU:HD21	1:A:1234:LEU:HB2	1.96	0.47
1:A:1298:PHE:CE1	1:A:1366:LEU:HD21	2.49	0.47
4:A:1601:FAD:H9	4:A:1601:FAD:C2'	2.45	0.47
1:A:1304:GLN:HE21	1:A:1373:GLY:HA3	1.79	0.47
1:A:294:GLU:HG2	1:A:557:PHE:CE2	2.49	0.47
1:A:98:LEU:HD23	1:A:556:VAL:HG11	1.97	0.46
1:A:1041:TYR:HE1	8:A:1608:PLC:H52	1.80	0.46
1:A:1133:ILE:HG21	1:A:1133:ILE:HD13	1.57	0.46
1:A:1235:LEU:HD21	6:A:1606:HEB:HBC1	1.98	0.46
1:A:1241:PHE:CD2	1:A:1243:LEU:HG	2.50	0.46
1:A:1323:LEU:HD22	1:A:1332:LEU:HD13	1.97	0.46
1:A:82:VAL:HG11	1:A:468:THR:HB	1.97	0.46
1:A:61:ASP:OD1	1:A:62:GLY:N	2.48	0.46
1:A:1455:THR:HG22	1:A:1495:ARG:HA	1.97	0.46
1:A:320:PRO:HB3	1:A:430:LEU:HD23	1.97	0.46
1:A:241:ARG:O	1:A:242:GLU:HG3	2.16	0.46
1:A:1113:ARG:O	1:A:1550:ASN:ND2	2.50	0.45
1:A:591:SER:HB2	1:A:1182:TYR:CE1	2.52	0.45
1:A:1113:ARG:NH2	1:A:1551:PHE:O	2.35	0.45
2:B:228:THR:HA	2:B:231:ILE:HG22	1.98	0.45
1:A:525:SER:OG	1:A:526:LYS:N	2.50	0.45
1:A:1287:VAL:HG11	1:A:1431:TRP:HE1	1.82	0.45
1:A:1499:GLU:HB2	1:A:1500:PRO:HD3	1.98	0.45
2:B:115:THR:HB	2:B:116:PRO:HD3	1.98	0.45
2:B:172:GLN:NE2	2:B:230:LEU:HB3	2.28	0.45
1:A:1225:HIS:HE1	6:A:1605:HEB:ND	2.15	0.45
1:A:338:VAL:HG11	1:A:388:MET:HG2	1.99	0.44
1:A:485:GLY:HA3	1:A:497:PHE:CZ	2.52	0.44
7:A:1607:NDP:C3D	8:A:1608:PLC:H71	2.47	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:56:ARG:HH21	2:B:198:LEU:HD11	1.83	0.44
1:A:332:THR:OG1	1:A:397:ASP:O	2.25	0.44
1:A:145:PRO:HG2	1:A:413:LYS:HB3	2.00	0.44
1:A:246:GLN:O	1:A:250:LEU:HG	2.17	0.44
1:A:553:GLN:HB3	1:A:563:ASP:HB3	2.00	0.43
1:A:1287:VAL:HG12	1:A:1337:ARG:HG2	1.99	0.43
1:A:1366:LEU:CD2	1:A:1368:GLY:H	2.31	0.43
1:A:1295:PRO:HG2	1:A:1298:PHE:HB2	1.99	0.43
1:A:1383:SER:HB3	1:A:1417:ILE:HG13	1.99	0.43
1:A:1377:TRP:NE1	1:A:1402:LEU:HD21	2.33	0.43
1:A:1526:THR:HG22	1:A:1548:TYR:CZ	2.53	0.43
2:B:156:TYR:CE1	2:B:160:LYS:HE2	2.53	0.43
1:A:1462:ASP:O	1:A:1464:ARG:N	2.51	0.43
1:A:76:ARG:O	1:A:80:ASN:ND2	2.51	0.43
1:A:125:ILE:HG12	1:A:425:GLN:HG3	2.01	0.43
1:A:508:LEU:HD23	1:A:508:LEU:HA	1.74	0.43
1:A:1132:LEU:O	1:A:1132:LEU:CG	2.67	0.43
1:A:516:PHE:CE2	1:A:517:GLU:HG3	2.53	0.42
1:A:1382:VAL:HB	1:A:1515:ILE:HG21	2.00	0.42
1:A:1495:ARG:HG3	1:A:1496:PRO:HD2	1.99	0.42
1:A:74:ASN:HB3	1:A:77:HIS:HB3	2.01	0.42
1:A:582:LEU:HD23	1:A:583:PHE:N	2.34	0.42
1:A:1090:ALA:HB2	1:A:1144:HIS:CE1	2.53	0.42
1:A:1235:LEU:HD23	1:A:1235:LEU:HA	1.87	0.42
8:A:1608:PLC:H83	8:A:1608:PLC:H42	1.75	0.42
1:A:121:GLU:HG3	1:A:148:ARG:HD3	2.00	0.42
1:A:150:ARG:HG3	1:A:168:GLN:NE2	2.35	0.42
1:A:1059:PHE:HD1	1:A:1060:LEU:HD12	1.83	0.42
1:A:158:SER:OG	1:A:160:SER:OG	2.34	0.42
1:A:1083:ILE:HD11	1:A:1244:ILE:HD13	2.02	0.42
1:A:98:LEU:HD23	1:A:556:VAL:CG1	2.49	0.42
1:A:1460:LYS:HD2	1:A:1463:LEU:HD11	2.01	0.42
1:A:518:ASN:HB3	1:A:521:ASN:HD21	1.85	0.41
7:A:1607:NDP:O3D	8:A:1608:PLC:H71	2.20	0.41
1:A:1415:LYS:HA	1:A:1415:LYS:HD2	1.95	0.41
1:A:297:PRO:HG3	1:A:304:PRO:HD3	2.02	0.41
7:A:1607:NDP:HO3N	8:A:1608:PLC:H81	1.79	0.41
1:A:559:TRP:CZ2	1:A:566:PRO:HA	2.56	0.41
2:B:183:MET:HB2	2:B:183:MET:HE2	1.85	0.41
1:A:149:SER:HB3	1:A:426:ARG:NH1	2.35	0.41
1:A:461:ASN:O	1:A:461:ASN:ND2	2.53	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1118:ASN:HD22	1:A:1550:ASN:ND2	2.18	0.41
1:A:1278:VAL:N	1:A:1291:ARG:O	2.47	0.41
1:A:110:LEU:HD21	1:A:173:LEU:HD13	2.02	0.41
1:A:1132:LEU:O	1:A:1132:LEU:CD1	2.69	0.41
1:A:1109:ILE:HG23	1:A:1123:PHE:HB3	2.01	0.41
1:A:93:ARG:NE	1:A:492:ASP:OD1	2.48	0.41
1:A:1230:PHE:CZ	1:A:1234:LEU:HD11	2.56	0.41
1:A:1318:TYR:C	1:A:1319:HIS:HD1	2.24	0.41
1:A:1524:GLY:O	7:A:1607:NDP:N6A	2.54	0.41
1:A:251:LEU:HD11	1:A:540:ILE:HG22	2.02	0.40
1:A:1130:HIS:HE1	6:A:1605:HEB:NC	2.19	0.40
1:A:1538:ARG:H	1:A:1538:ARG:HG2	1.75	0.40
1:A:145:PRO:HB2	1:A:412:LEU:HD21	2.04	0.40
1:A:1045:ILE:HG23	8:A:1608:PLC:H6'2	2.03	0.40
1:A:286:THR:HG21	1:A:540:ILE:HD11	2.04	0.40
2:B:63:ILE:O	2:B:66:VAL:HG12	2.21	0.40
1:A:67:LEU:HD22	1:A:71:TYR:HD2	1.85	0.40
2:B:232:SER:HB3	2:B:233:PRO:HD3	2.02	0.40
1:A:482:LEU:HB3	1:A:483:PRO:HD3	2.03	0.40
1:A:1062:ARG:HH21	1:A:1248:ARG:NH2	2.19	0.40
1:A:1108:LEU:HB2	1:A:1263:ASP:OD2	2.22	0.40
1:A:1133:ILE:O	1:A:1133:ILE:CG2	2.69	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	1075/1536 (70%)	1012 (94%)	61 (6%)	2 (0%)	44	71
2	B	226/341 (66%)	216 (96%)	10 (4%)	0	100	100
All	All	1301/1877 (69%)	1228 (94%)	71 (6%)	2 (0%)	45	71

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	1352	PRO
1	A	1463	LEU

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	879/1348 (65%)	874 (99%)	5 (1%)	84	90
2	B	179/287 (62%)	179 (100%)	0	100	100
All	All	1058/1635 (65%)	1053 (100%)	5 (0%)	85	91

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	211	SER
1	A	213	LEU
1	A	214	LEU
1	A	1463	LEU
1	A	1464	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	289	ASN
1	A	348	GLN
1	A	451	GLN
1	A	504	GLN
1	A	1107	ASN
1	A	1130	HIS
1	A	1226	HIS
1	A	1238	HIS
1	A	1245	GLN
1	A	1375	GLN

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Mol	Chain	Res	Type
1	A	1537	ASN
1	A	1550	ASN
2	B	71	ASN
2	B	172	GLN

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 ⓘ

5 monosaccharides 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	$\# Z > 2$
3	NAG	C	1	3,2	14,14,15	0.24	0	17,19,21	0.56	0
3	NAG	C	2	3	14,14,15	0.36	0	17,19,21	0.34	0
3	BMA	C	3	3	11,11,12	0.59	0	15,15,17	0.86	0
3	MAN	C	4	3	11,11,12	1.00	0	15,15,17	0.88	0
3	MAN	C	5	3	11,11,12	0.63	0	15,15,17	1.19	2 (13%)

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
3	NAG	C	1	3,2	-	0/6/23/26	0/1/1/1
3	NAG	C	2	3	-	1/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	BMA	C	3	3	-	1/2/19/22	0/1/1/1
3	MAN	C	4	3	-	2/2/19/22	0/1/1/1
3	MAN	C	5	3	-	2/2/19/22	0/1/1/1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	5	MAN	C1-O5-C5	2.99	116.19	112.19
3	C	5	MAN	O2-C2-C3	-2.29	105.40	110.15

There are no chirality outliers.

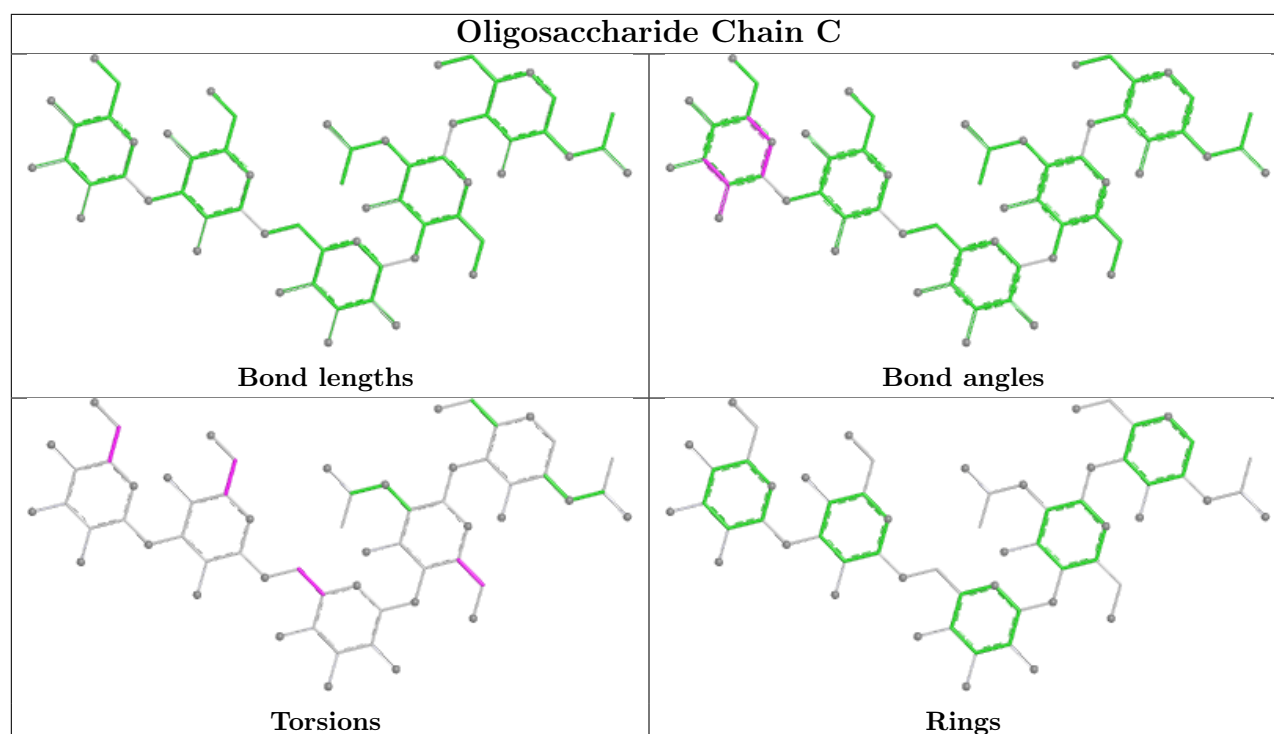
All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	C	4	MAN	O5-C5-C6-O6
3	C	4	MAN	C4-C5-C6-O6
3	C	5	MAN	O5-C5-C6-O6
3	C	2	NAG	O5-C5-C6-O6
3	C	5	MAN	C4-C5-C6-O6
3	C	3	BMA	O5-C5-C6-O6

There are no ring outliers.

No monomer is involved in short contacts.

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.



5.6 Ligand geometry [i](#)

Of 10 ligands modelled in this entry, 1 is monoatomic - leaving 9 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	$\# Z > 2$
8	PLC	A	1608	-	41,41,41	1.06	2 (4%)	47,49,49	1.21	4 (8%)
4	FAD	A	1601	-	54,58,58	1.36	7 (12%)	71,89,89	1.50	14 (19%)
5	NAG	A	1602	1	14,14,15	0.22	0	17,19,21	0.48	0
7	NDP	A	1607	-	38,43,52	1.28	3 (7%)	52,67,80	2.00	10 (19%)
5	NAG	A	1603	1	14,14,15	0.40	0	17,19,21	1.37	2 (11%)
5	NAG	A	1604	1	14,14,15	1.24	1 (7%)	17,19,21	1.41	1 (5%)
6	HEB	A	1606	1	47,50,50	1.72	12 (25%)	55,82,82	2.02	17 (30%)
6	HEB	A	1605	1	47,50,50	1.70	12 (25%)	55,82,82	2.10	20 (36%)
5	NAG	B	406	2	14,14,15	0.24	0	17,19,21	0.52	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
8	PLC	A	1608	-	-	13/45/45/45	-
4	FAD	A	1601	-	-	15/30/50/50	0/6/6/6
7	NDP	A	1607	-	-	11/23/59/77	0/4/4/5
5	NAG	A	1603	1	-	6/6/23/26	0/1/1/1
5	NAG	A	1604	1	-	4/6/23/26	0/1/1/1
6	HEB	A	1606	1	1/1/3/8	2/12/54/54	-
5	NAG	B	406	2	-	2/6/23/26	0/1/1/1
6	HEB	A	1605	1	1/1/3/8	4/12/54/54	-
5	NAG	A	1602	1	-	2/6/23/26	0/1/1/1

All (37) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	1605	HEB	CBB-CAB	-4.63	1.31	1.51
4	A	1601	FAD	C4-N3	-4.51	1.30	1.38
6	A	1606	HEB	CBB-CAB	-4.45	1.32	1.51
5	A	1604	NAG	O5-C1	4.24	1.50	1.43
8	A	1608	PLC	O3-CB	4.16	1.45	1.33
8	A	1608	PLC	O2-C'	4.16	1.46	1.34
6	A	1606	HEB	C4B-C3B	-3.70	1.38	1.45
4	A	1601	FAD	C9A-C5X	3.68	1.47	1.41
7	A	1607	NDP	PN-O3	-3.57	1.55	1.59
6	A	1605	HEB	C4B-C3B	-3.55	1.39	1.45
6	A	1605	HEB	C4A-C3A	-3.55	1.38	1.45
6	A	1606	HEB	C1A-C2A	-3.36	1.39	1.45
6	A	1605	HEB	C4D-C3D	-3.34	1.39	1.45
6	A	1606	HEB	C4D-C3D	-3.33	1.39	1.45
4	A	1601	FAD	C2-N3	-3.31	1.31	1.39
6	A	1606	HEB	C4A-C3A	-3.10	1.39	1.45
6	A	1606	HEB	C4B-NB	-2.90	1.35	1.40
6	A	1605	HEB	C1D-ND	-2.89	1.35	1.40
7	A	1607	NDP	C3B-C4B	-2.87	1.45	1.53
6	A	1606	HEB	C1D-ND	-2.84	1.35	1.40
6	A	1605	HEB	C4B-NB	-2.75	1.35	1.40
6	A	1605	HEB	C1D-C2D	-2.74	1.39	1.44
6	A	1606	HEB	C1D-C2D	-2.71	1.39	1.44
6	A	1605	HEB	C1A-C2A	-2.67	1.40	1.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1601	FAD	C8-C7	2.60	1.47	1.40
4	A	1601	FAD	C5X-N5	-2.57	1.34	1.39
6	A	1606	HEB	C1B-C2B	-2.54	1.39	1.44
7	A	1607	NDP	O4B-C4B	-2.48	1.39	1.45
6	A	1606	HEB	C4A-NA	-2.48	1.35	1.39
6	A	1605	HEB	C1B-C2B	-2.46	1.39	1.44
6	A	1605	HEB	C4A-NA	-2.44	1.35	1.39
6	A	1606	HEB	C1A-NA	-2.35	1.35	1.39
4	A	1601	FAD	C6-C7	-2.25	1.36	1.39
6	A	1605	HEB	C1A-NA	-2.11	1.35	1.39
6	A	1605	HEB	CHC-C4B	2.04	1.39	1.34
6	A	1606	HEB	CHD-C1D	2.02	1.39	1.34
4	A	1601	FAD	C5A-N7A	-2.02	1.32	1.39

All (68) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	A	1607	NDP	O4B-C1B-C2B	-6.55	95.43	106.61
7	A	1607	NDP	N3A-C2A-N1A	-6.19	120.27	128.67
5	A	1604	NAG	C1-O5-C5	5.60	119.69	112.19
6	A	1605	HEB	CHB-C4A-NA	5.45	130.39	124.45
6	A	1605	HEB	CHA-C4D-ND	4.72	129.50	124.42
8	A	1608	PLC	O2-C'-C1'	4.63	121.49	111.48
5	A	1603	NAG	C2-N2-C7	4.59	129.05	122.90
6	A	1605	HEB	CHB-C1B-NB	4.52	129.28	124.42
6	A	1606	HEB	CHA-C4D-ND	4.41	129.16	124.42
6	A	1606	HEB	CHB-C1B-NB	4.34	129.09	124.42
6	A	1606	HEB	CHA-C1A-NA	4.32	129.16	124.45
4	A	1601	FAD	N3A-C2A-N1A	-4.31	122.82	128.67
6	A	1606	HEB	CHC-C4B-NB	4.22	129.60	124.37
6	A	1606	HEB	CHB-C4A-NA	4.13	128.95	124.45
6	A	1606	HEB	CHD-C1D-ND	4.11	129.47	124.37
7	A	1607	NDP	O3-PN-O1N	-4.00	98.67	110.70
6	A	1605	HEB	CHD-C1D-ND	3.96	129.28	124.37
6	A	1605	HEB	CHC-C4B-NB	3.94	129.25	124.37
7	A	1607	NDP	O3B-C3B-C4B	-3.80	100.16	111.08
6	A	1605	HEB	CBB-CAB-C3B	3.54	122.02	112.42
6	A	1606	HEB	CBB-CAB-C3B	3.52	121.95	112.42
6	A	1605	HEB	CAA-C2A-C1A	3.35	131.66	124.85
4	A	1601	FAD	C4X-C10-N1	-3.33	116.43	124.59
8	A	1608	PLC	C2-O2-C'	-3.31	109.89	117.80
4	A	1601	FAD	C4-C4X-N5	3.24	122.69	118.21

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1601	FAD	C4-N3-C2	-3.20	119.96	125.64
7	A	1607	NDP	O4D-C4D-C5D	3.20	119.58	109.33
7	A	1607	NDP	O2N-PN-O1N	3.17	127.19	112.44
6	A	1606	HEB	C4A-CHB-C1B	-3.16	119.30	126.02
4	A	1601	FAD	O4B-C1B-N9A	3.11	112.86	108.75
7	A	1607	NDP	O5B-C5B-C4B	-3.04	98.64	108.99
6	A	1605	HEB	CAA-C2A-C3A	-2.99	122.27	127.87
6	A	1605	HEB	CHB-C1B-C2B	-2.94	120.39	125.03
4	A	1601	FAD	C4X-C4-N3	2.93	120.71	113.25
6	A	1605	HEB	C4A-CHB-C1B	-2.92	119.80	126.02
6	A	1606	HEB	CHB-C1B-C2B	-2.90	120.46	125.03
4	A	1601	FAD	C4X-C10-N10	2.89	120.62	116.48
7	A	1607	NDP	C5B-C4B-C3B	-2.89	104.82	115.21
6	A	1606	HEB	C1A-CHA-C4D	-2.84	119.98	126.02
6	A	1605	HEB	C1A-CHA-C4D	-2.76	120.15	126.02
8	A	1608	PLC	O3-CB-C1B	2.71	120.11	111.83
6	A	1605	HEB	CHA-C4D-C3D	-2.60	120.98	124.77
6	A	1605	HEB	CHB-C4A-C3A	-2.57	120.15	125.49
7	A	1607	NDP	C4D-O4D-C1D	-2.54	103.35	108.63
6	A	1605	HEB	CHA-C1A-NA	2.47	127.14	124.45
7	A	1607	NDP	O2B-P2B-O1X	-2.45	100.60	109.33
4	A	1601	FAD	C10-N1-C2	2.44	122.12	116.85
4	A	1601	FAD	C4'-C3'-C2'	-2.42	109.55	113.57
6	A	1606	HEB	CAD-CBD-CGD	-2.37	107.38	113.67
6	A	1605	HEB	C4B-NB-C1B	2.37	108.01	105.21
6	A	1606	HEB	C1B-C2B-C3B	-2.28	104.59	106.98
6	A	1606	HEB	C4B-NB-C1B	2.26	107.88	105.21
6	A	1606	HEB	C1D-ND-C4D	2.24	107.86	105.21
6	A	1606	HEB	CHA-C1A-C2A	-2.22	121.35	124.86
6	A	1606	HEB	CHA-C4D-C3D	-2.22	121.53	124.77
5	A	1603	NAG	C1-C2-N2	2.20	113.91	110.43
6	A	1605	HEB	CAA-CBA-CGA	-2.20	107.82	113.67
4	A	1601	FAD	C9A-N10-C10	-2.16	117.46	120.75
6	A	1606	HEB	CHD-C1D-C2D	-2.15	120.85	126.94
6	A	1605	HEB	C1B-C2B-C3B	-2.12	104.75	106.98
4	A	1601	FAD	C4A-C5A-N7A	-2.12	107.10	109.34
6	A	1605	HEB	CHD-C1D-C2D	-2.12	120.94	126.94
6	A	1605	HEB	C1D-ND-C4D	2.11	107.71	105.21
4	A	1601	FAD	N3-C2-N1	2.11	123.98	119.50
4	A	1601	FAD	C10-C4X-N5	-2.10	120.52	124.81
8	A	1608	PLC	O2-C'-O'	-2.09	118.81	123.70
4	A	1601	FAD	O4'-C4'-C3'	-2.01	104.55	109.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	1605	HEB	CMC-C2C-C3C	2.01	128.70	124.68

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
6	A	1605	HEB	NA
6	A	1606	HEB	NA

All (59) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	1601	FAD	C5B-O5B-PA-O1A
4	A	1601	FAD	C5B-O5B-PA-O3P
4	A	1601	FAD	N10-C1'-C2'-O2'
4	A	1601	FAD	N10-C1'-C2'-C3'
4	A	1601	FAD	C5'-O5'-P-O2P
7	A	1607	NDP	C5B-O5B-PA-O1A
7	A	1607	NDP	C5B-O5B-PA-O2A
7	A	1607	NDP	C5B-O5B-PA-O3
7	A	1607	NDP	C5D-O5D-PN-O3
7	A	1607	NDP	O4D-C4D-C5D-O5D
8	A	1608	PLC	O4P-C4-C5-N
8	A	1608	PLC	C1-O3P-P-O1P
8	A	1608	PLC	C1-O3P-P-O2P
8	A	1608	PLC	C1-O3P-P-O4P
6	A	1605	HEB	C1A-C2A-CAA-CBA
6	A	1605	HEB	C2B-C3B-CAB-CBB
6	A	1606	HEB	C2B-C3B-CAB-CBB
5	B	406	NAG	O5-C5-C6-O6
5	A	1604	NAG	C4-C5-C6-O6
5	A	1602	NAG	O5-C5-C6-O6
6	A	1605	HEB	C3A-C2A-CAA-CBA
5	A	1604	NAG	O5-C5-C6-O6
5	A	1602	NAG	C4-C5-C6-O6
5	A	1603	NAG	C4-C5-C6-O6
6	A	1605	HEB	C4B-C3B-CAB-CBB
6	A	1606	HEB	C4B-C3B-CAB-CBB
5	B	406	NAG	C4-C5-C6-O6
5	A	1603	NAG	C8-C7-N2-C2
5	A	1603	NAG	O7-C7-N2-C2
5	A	1604	NAG	C8-C7-N2-C2
5	A	1604	NAG	O7-C7-N2-C2

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Mol	Chain	Res	Type	Atoms
7	A	1607	NDP	C3D-C4D-C5D-O5D
5	A	1603	NAG	O5-C5-C6-O6
7	A	1607	NDP	O4B-C4B-C5B-O5B
7	A	1607	NDP	C3B-C4B-C5B-O5B
8	A	1608	PLC	C1'-C'-O2-C2
8	A	1608	PLC	O'-C'-O2-C2
4	A	1601	FAD	C1'-C2'-C3'-O3'
8	A	1608	PLC	O3P-C1-C2-O2
8	A	1608	PLC	O2-C2-C3-O3
4	A	1601	FAD	C5B-O5B-PA-O2A
4	A	1601	FAD	C2'-C1'-N10-C10
4	A	1601	FAD	C5'-O5'-P-O3P
7	A	1607	NDP	C5D-O5D-PN-O1N
8	A	1608	PLC	O3P-C1-C2-C3
7	A	1607	NDP	PN-O3-PA-O1A
7	A	1607	NDP	PN-O3-PA-O2A
8	A	1608	PLC	C'-C1'-C2'-C3'
4	A	1601	FAD	O2'-C2'-C3'-O3'
4	A	1601	FAD	PA-O3P-P-O1P
5	A	1603	NAG	C1-C2-N2-C7
5	A	1603	NAG	C3-C2-N2-C7
4	A	1601	FAD	O2'-C2'-C3'-C4'
8	A	1608	PLC	C4B-C5B-C6B-C7B
8	A	1608	PLC	C1-C2-C3-O3
4	A	1601	FAD	O4B-C4B-C5B-O5B
8	A	1608	PLC	C2B-C3B-C4B-C5B
4	A	1601	FAD	P-O3P-PA-O2A
4	A	1601	FAD	PA-O3P-P-O2P

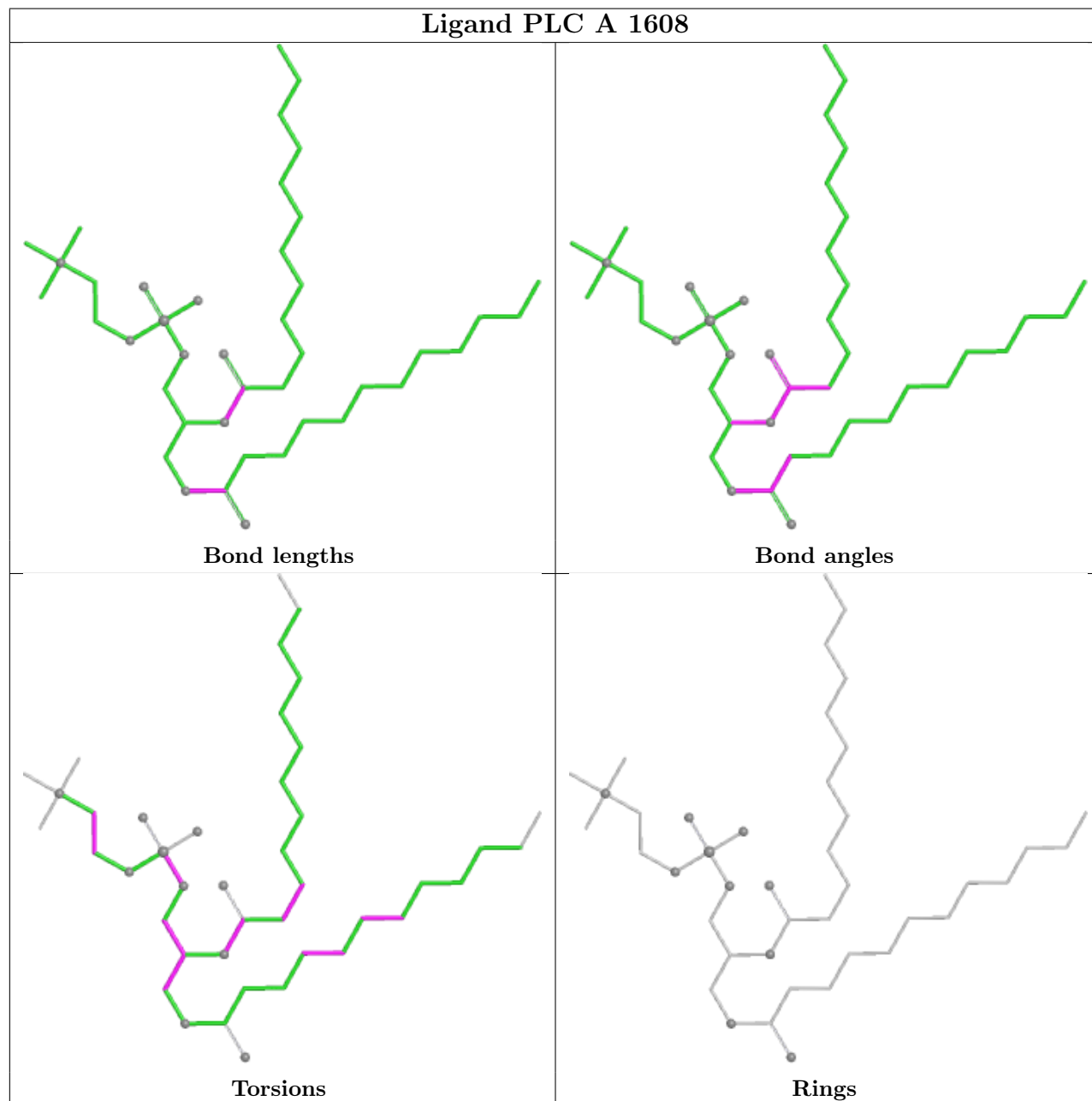
There are no ring outliers.

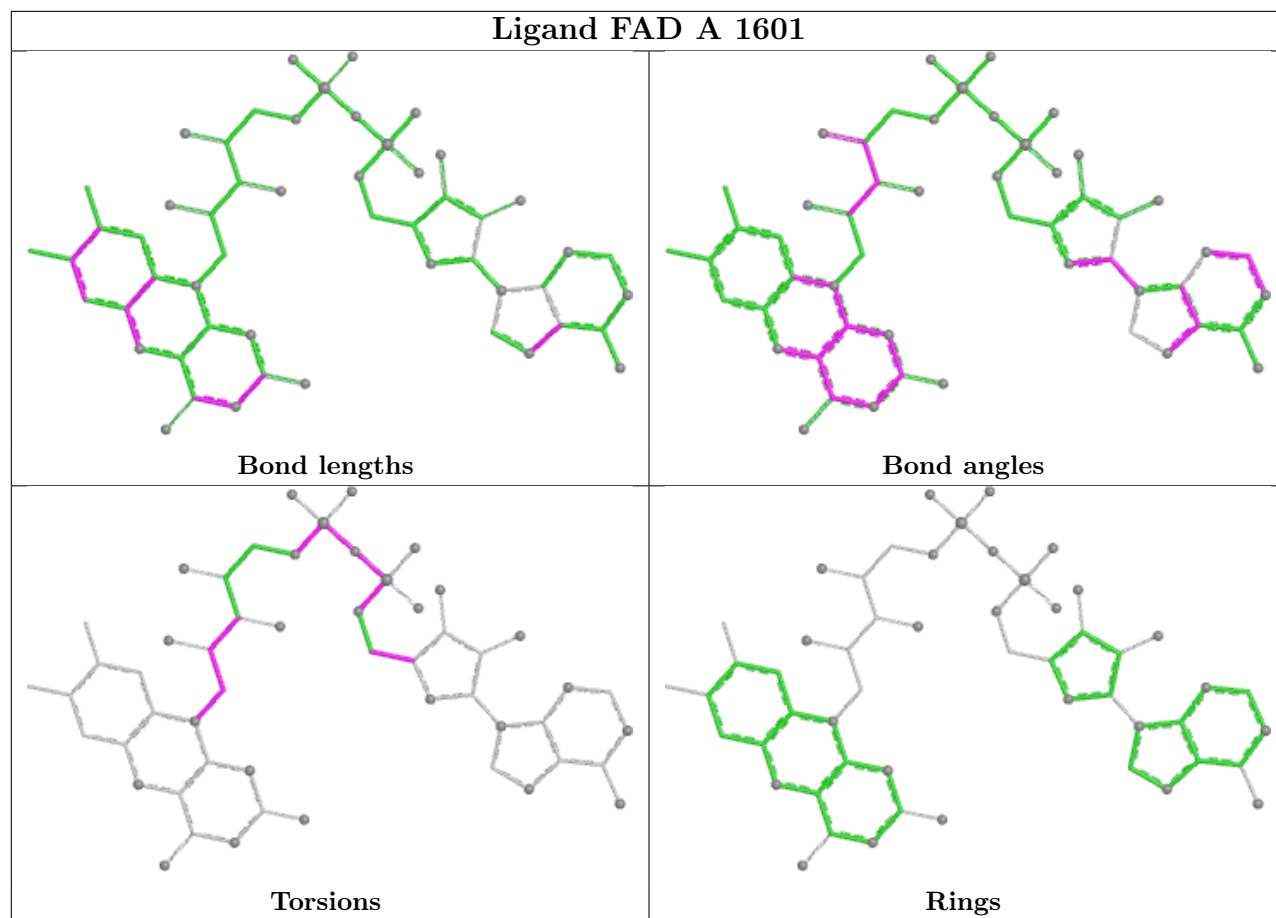
6 monomers are involved in 23 short contacts:

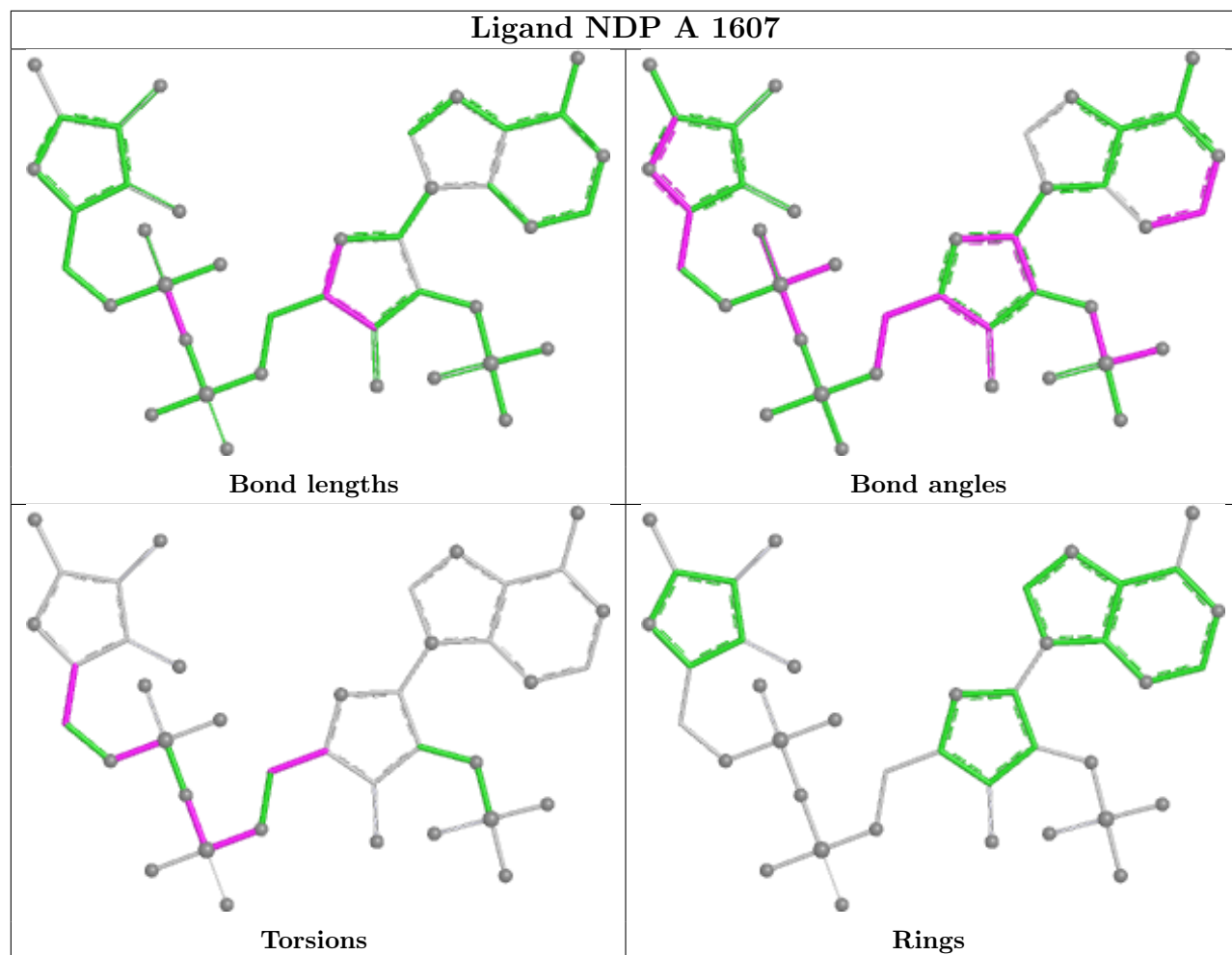
Mol	Chain	Res	Type	Clashes	Symm-Clashes
8	A	1608	PLC	11	0
4	A	1601	FAD	3	0
7	A	1607	NDP	8	0
5	A	1603	NAG	1	0
6	A	1606	HEB	3	0
6	A	1605	HEB	4	0

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

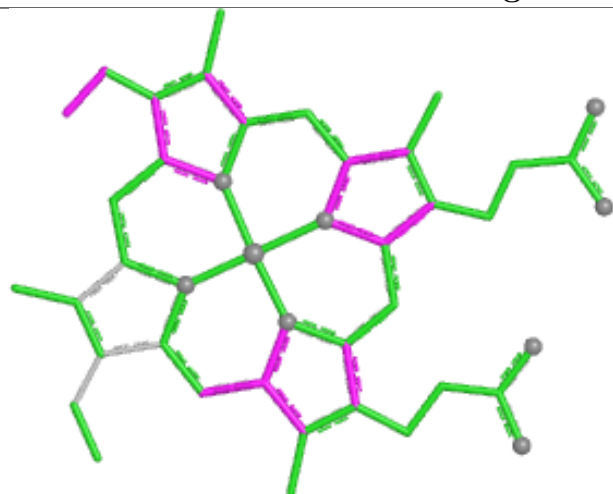
addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. 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.



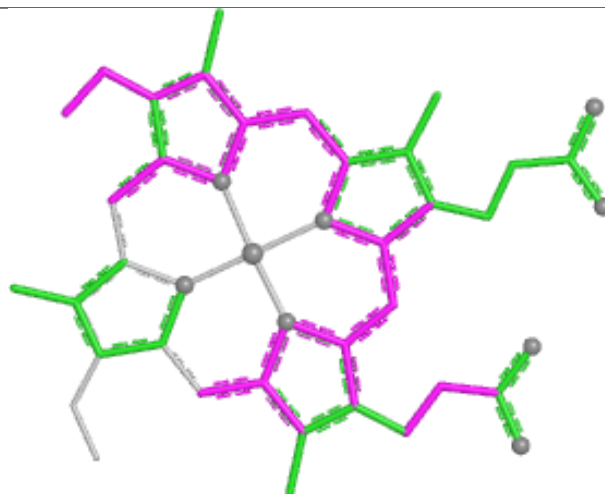




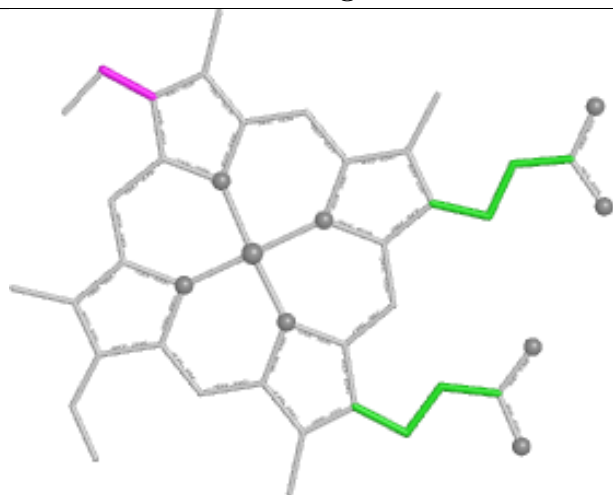
Ligand HEB A 1606



Bond lengths



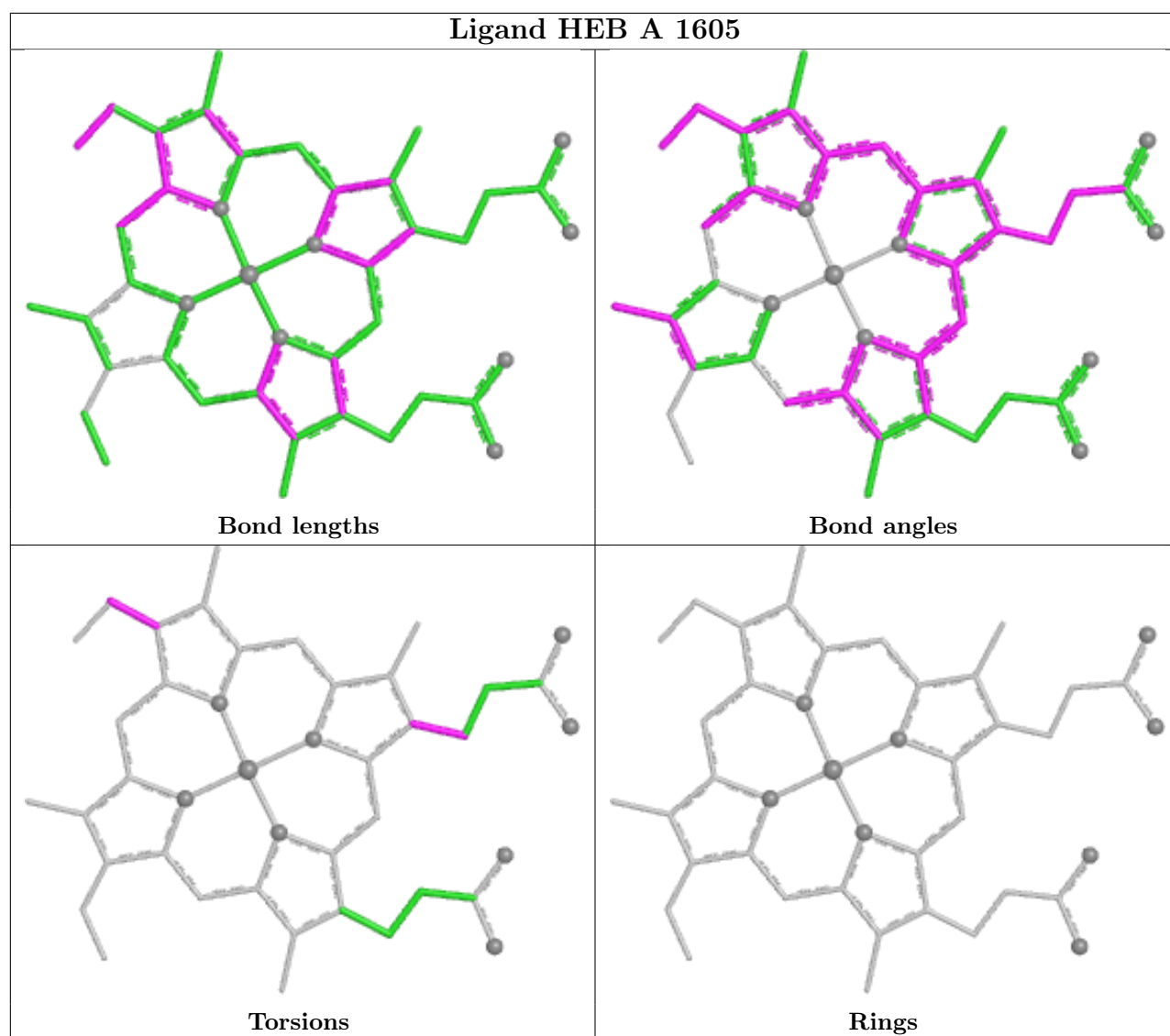
Bond angles



Torsions



Rings



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.

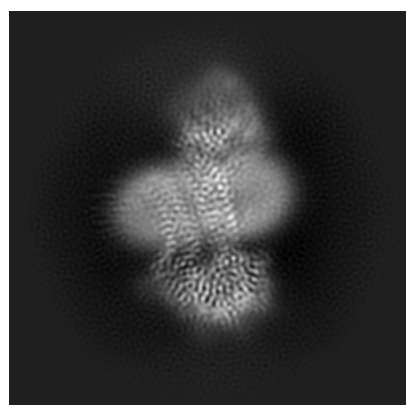
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-21964. These allow visual inspection of the internal detail of the map and identification of artifacts.

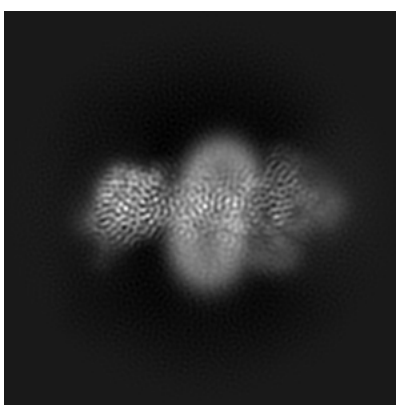
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

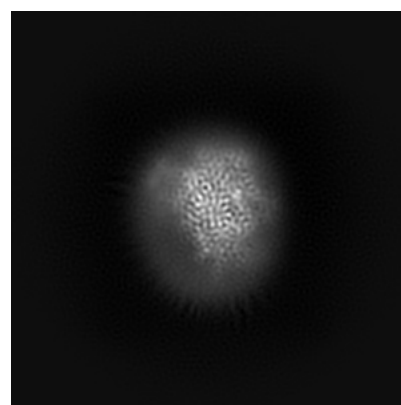
6.1.1 Primary map



X



Y

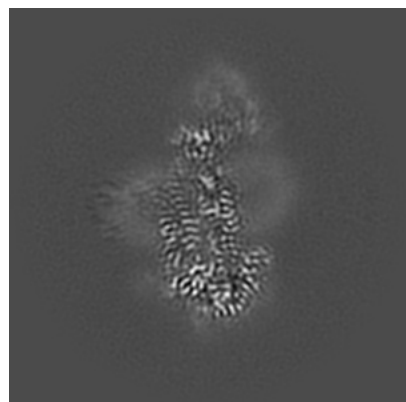


Z

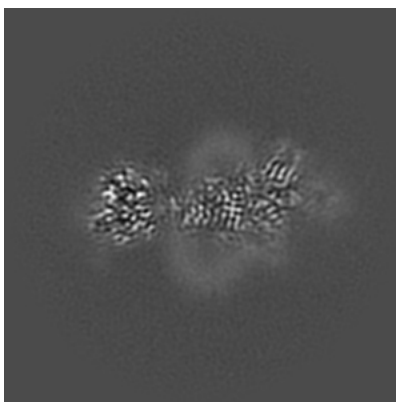
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

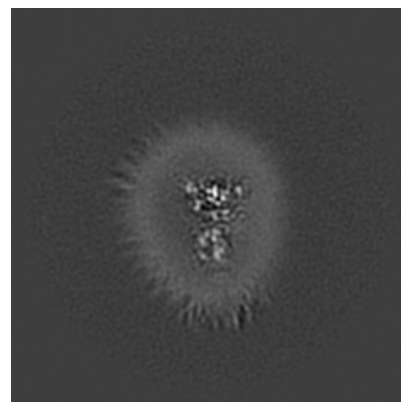
6.2.1 Primary map



X Index: 120



Y Index: 120

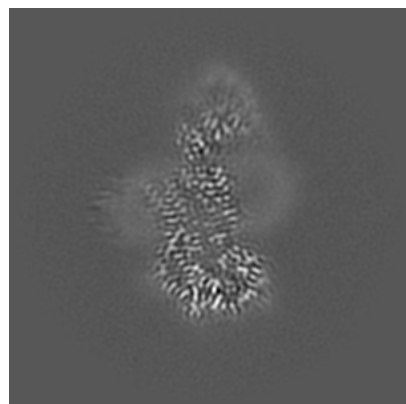


Z Index: 120

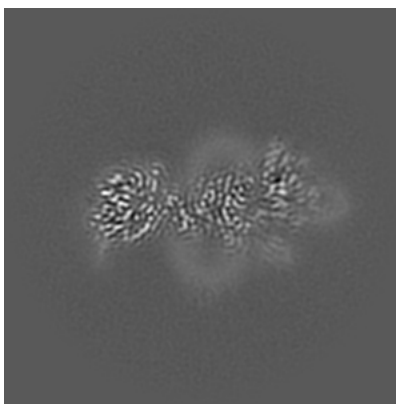
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

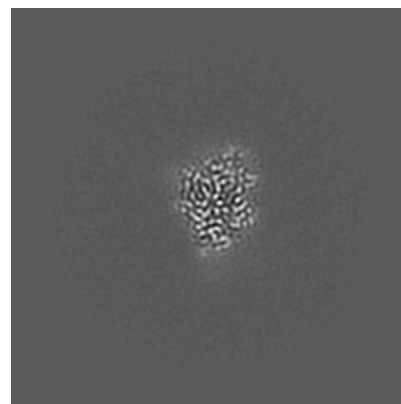
6.3.1 Primary map



X Index: 125



Y Index: 126

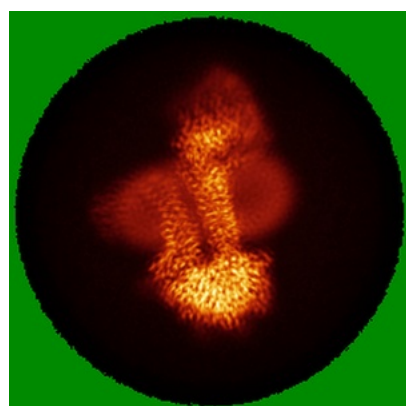


Z Index: 70

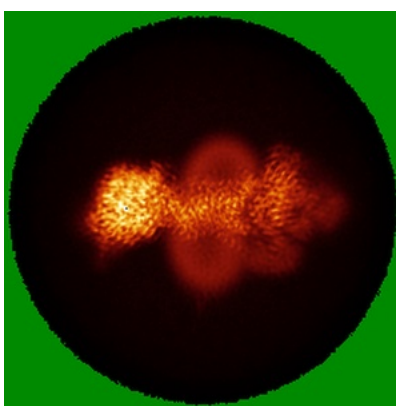
The images above show the largest variance slices of the map in three orthogonal directions.

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

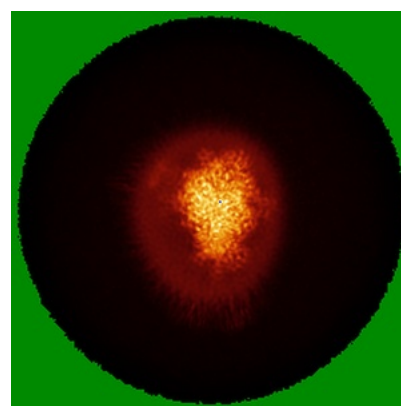
6.4.1 Primary map



X



Y

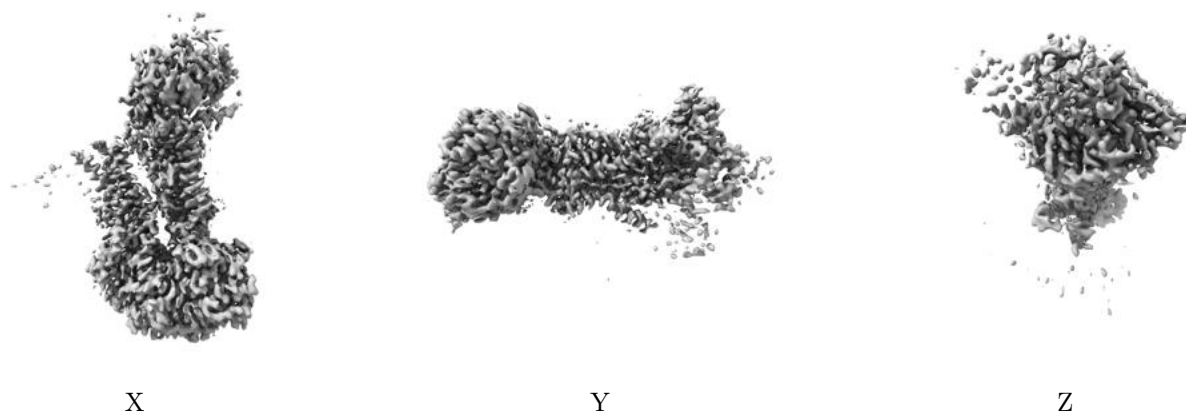


Z

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 0.55. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

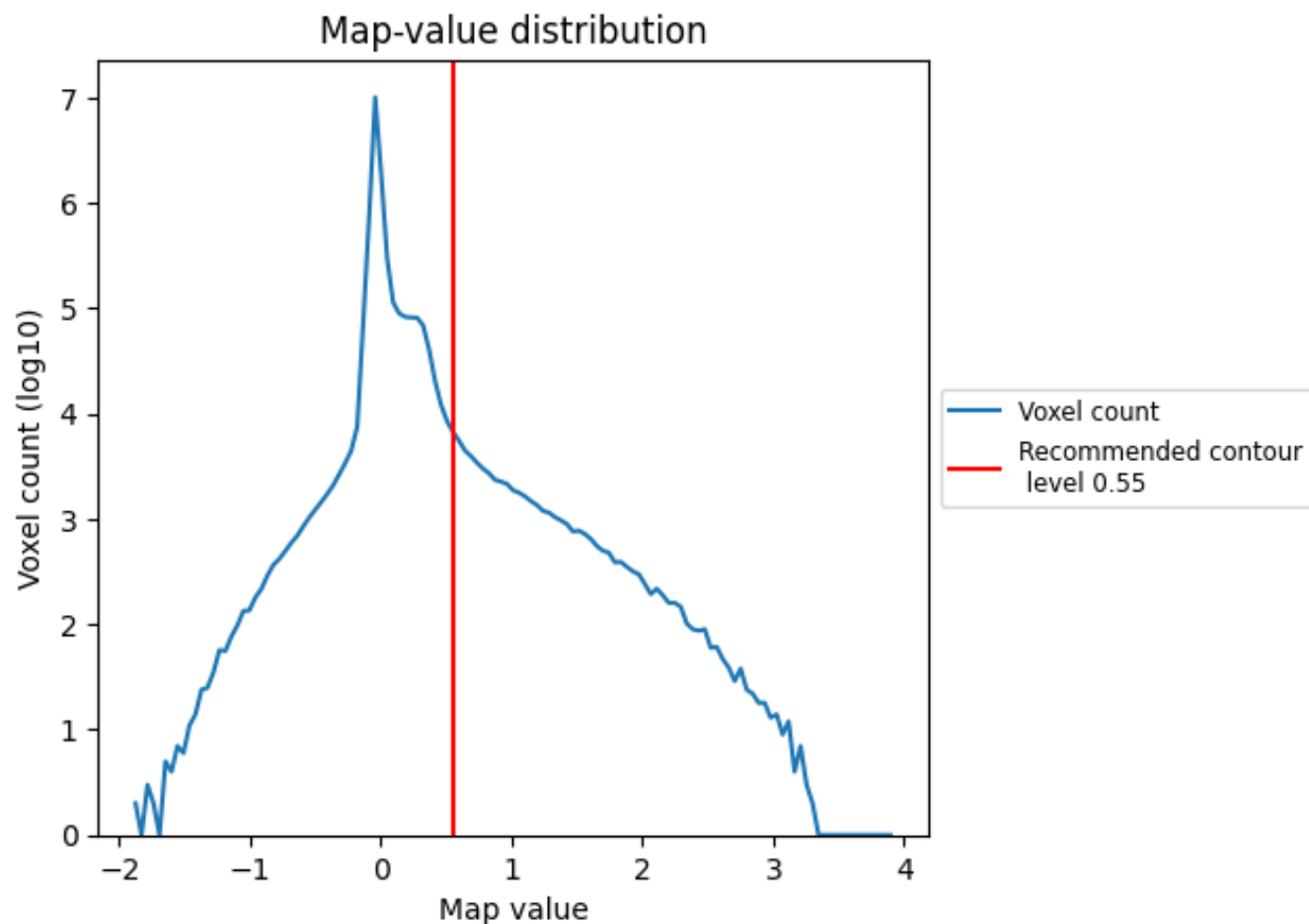
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

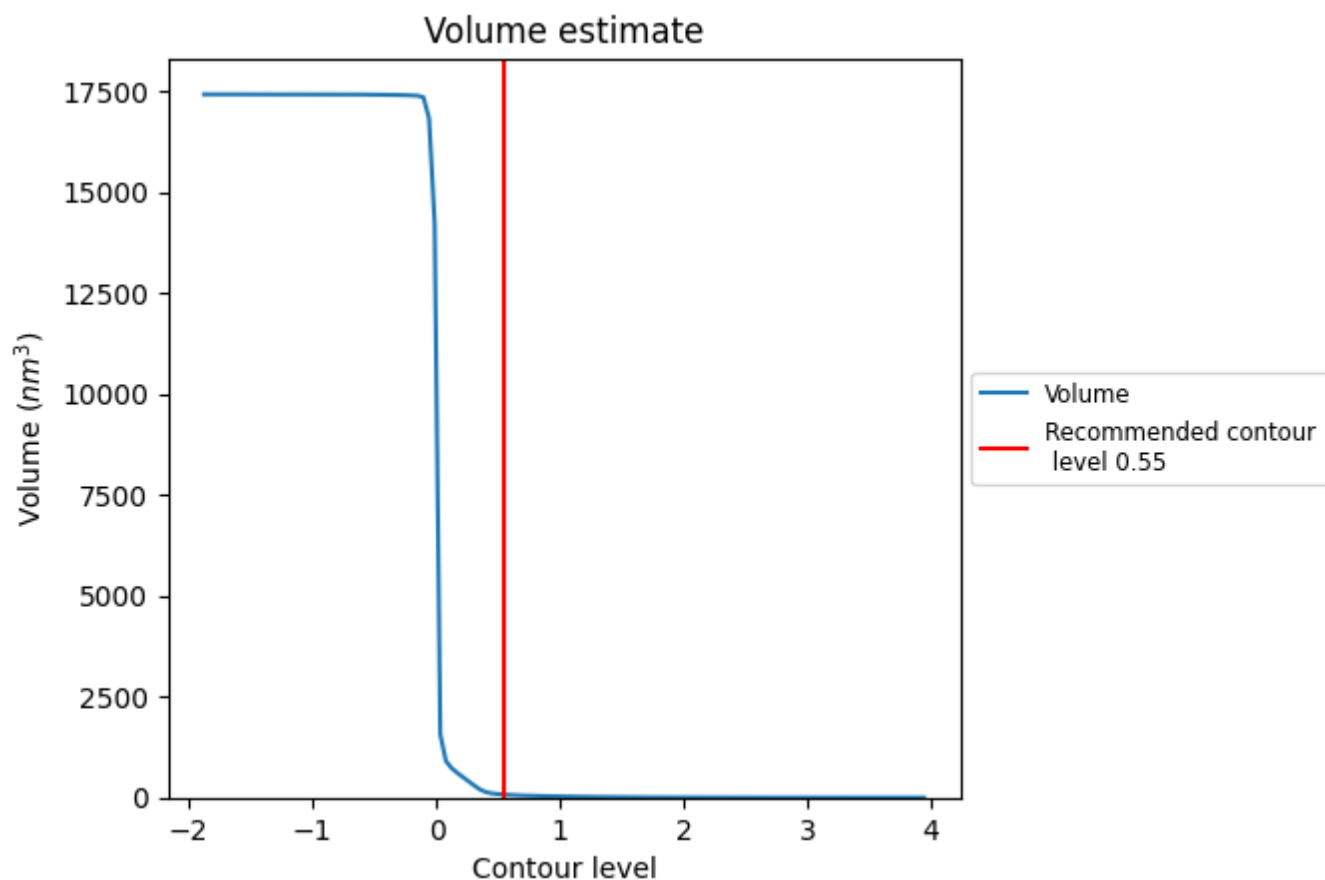
This section contains the results of statistical analysis of the map.

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.

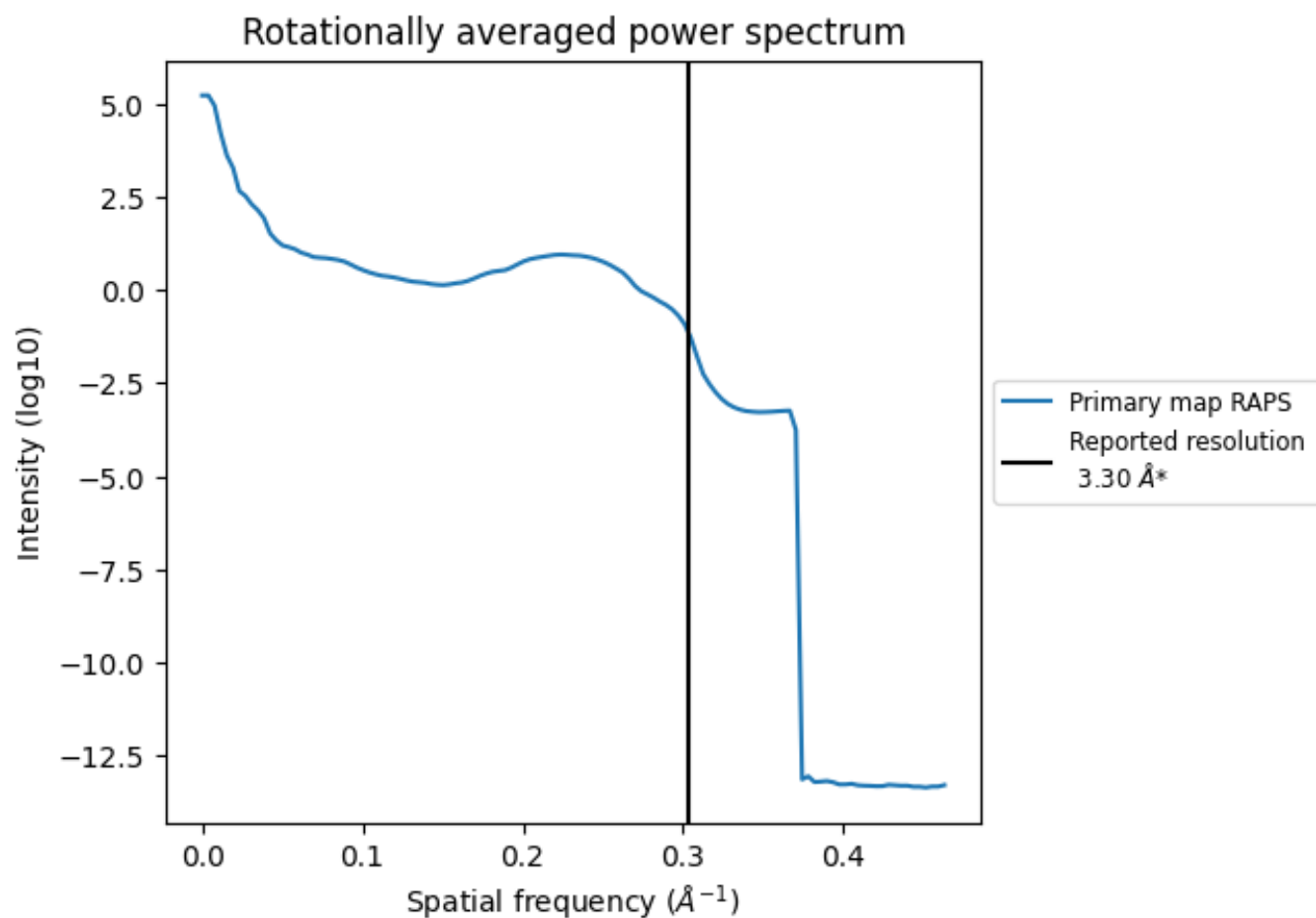
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 74 nm³; this corresponds to an approximate mass of 67 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

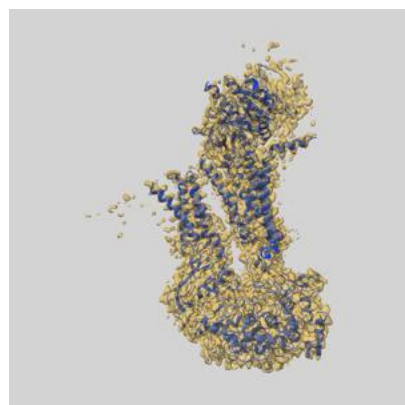
8 Fourier-Shell correlation

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

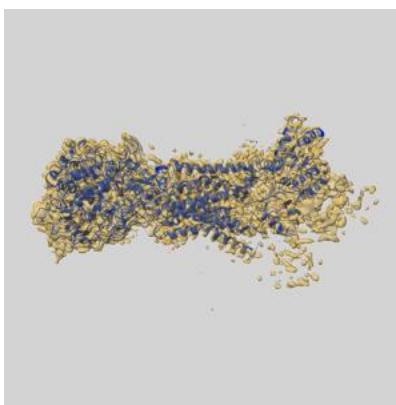
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-21964 and PDB model 6WXV. Per-residue inclusion information can be found in section [3](#) on page [8](#).

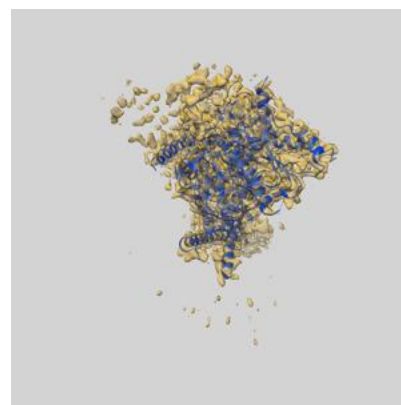
9.1 Map-model overlay [i](#)



X



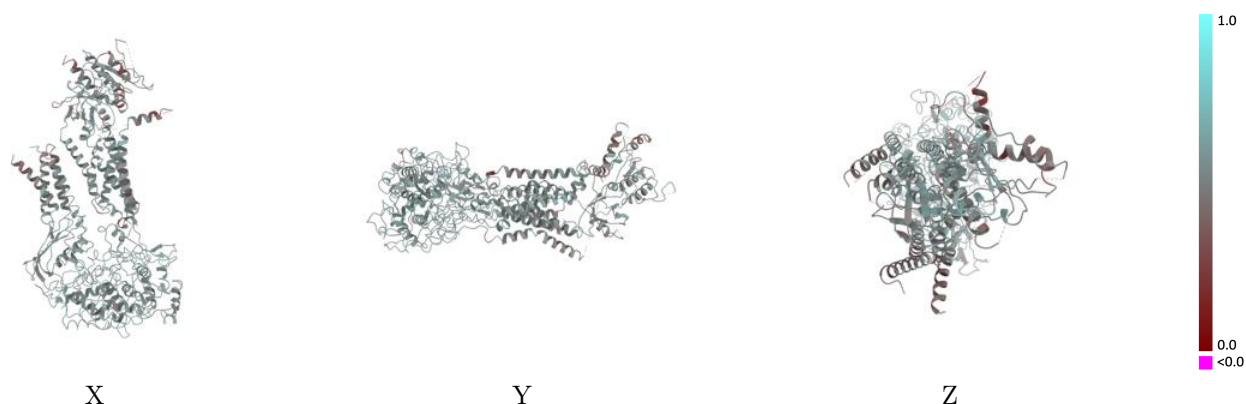
Y



Z

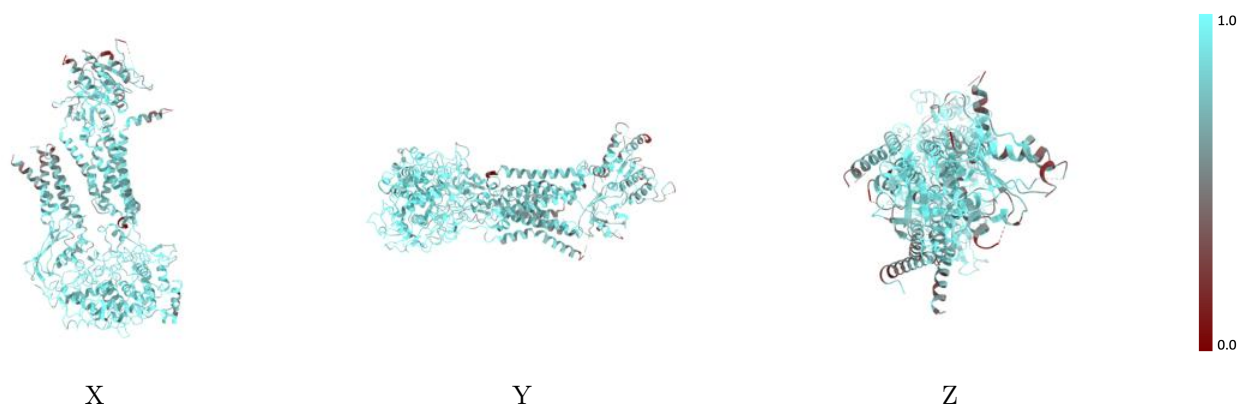
The images above show the 3D surface view of the map at the recommended contour level 0.55 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



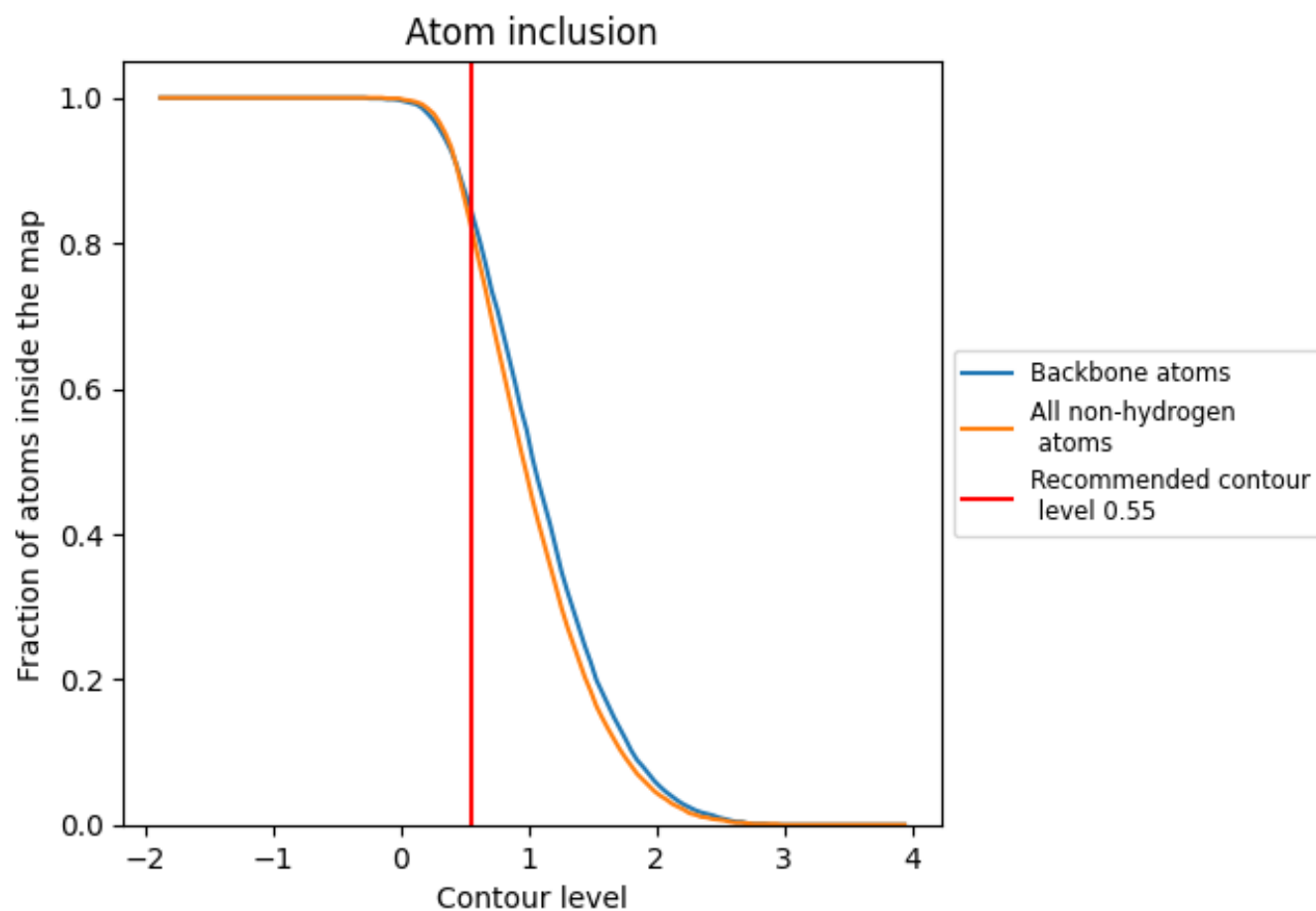
The images above show the model with each residue coloured according its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.55).

9.4 Atom inclusion [i](#)



At the recommended contour level, 85% 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 (0.55) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	<div></div> 0.8260	<div></div> 0.5260
A	<div></div> 0.8410	<div></div> 0.5300
B	<div></div> 0.7570	<div></div> 0.5070
C	<div></div> 0.6720	<div></div> 0.4540

