Full wwPDB EM Validation Report

Nov 4, 2023 – 03:54 PM EDT

PDB ID	:	8G5F
EMDB ID	:	EMD-29741
Title	:	Native GABA-A receptor from the mouse brain, ortho-alpha1-alpha3-beta2-gamma2 subtype, in complex with GABA and allopregnanolone
Authors	:	Sun, C.; Gouaux, E.
Deposited on	:	2023-02-13
Resolution	:	2.64 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:

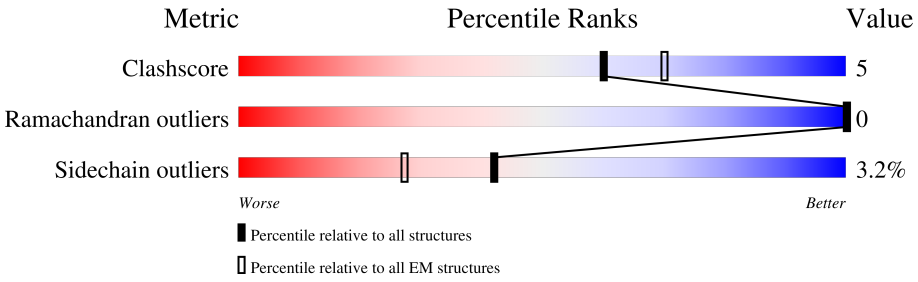
EMDB validation analysis	:	0.0.1.dev70
Mogul	:	1.8.5 (274361), CSD as541be (2020)
MolProbity	:	4.02b-467
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
MapQ	:	1.9.9
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

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 2.64 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	158937	4297
Ramachandran outliers	154571	4023
Sidechain outliers	154315	3826

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	492	59% 10% 30%
2	B	512	54% 10% 35%
2	E	512	53% 11% 35%
3	C	455	66% 9% 25%
4	D	474	5% 56% 10% 33%
5	J	223	46% 6% 48%
6	K	213	45% 51%
7	F	7	29% 71%

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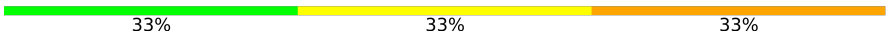
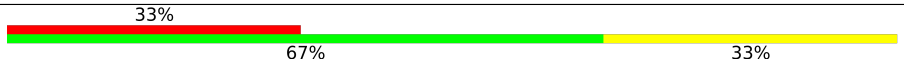
Mol	Chain	Length	Quality of chain
8	G	3	
8	H	3	
8	M	3	
8	N	3	
9	I	2	
9	L	2	

2 Entry composition (i)

There are 12 unique types of molecules in this entry. The entry contains 31366 atoms, of which 15604 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 Gamma-aminobutyric acid receptor subunit alpha-3.

Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	S		
1	A	343	5544	1807	2768	458	497	14	0	0

- Molecule 2 is a protein called Gamma-aminobutyric acid receptor subunit beta-2.

Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	S		
2	B	331	5429	1779	2717	434	483	16	0	0
2	E	331	Total	C	H	N	O	S	0	0
			5430	1779	2716	435	484	16		

- Molecule 3 is a protein called Gamma-aminobutyric acid receptor subunit alpha-1.

Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	S		
3	C	343	5565	1809	2774	468	498	16	1	0

- Molecule 4 is a protein called Gamma-aminobutyric acid receptor subunit gamma-2.

Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	S		
4	D	318	5187	1698	2582	428	465	14	0	0

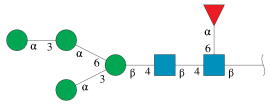
- Molecule 5 is a protein called Heavy Chain of 8E3 Fab.

Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	S		
5	J	117	1783	587	863	152	177	4	0	0

- Molecule 6 is a protein called Light Chain of 8E3 Fab.

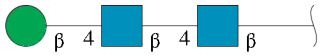
Mol	Chain	Residues	Atoms						AltConf	Trace
6	K	105	Total	C	H	N	O	S	0	0
			1552	505	756	132	155	4		

- Molecule 7 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms					AltConf	Trace
7	F	7	Total	C	H	N	O	0	0
			152	46	70	2	34		

- Molecule 8 is an oligosaccharide called beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.

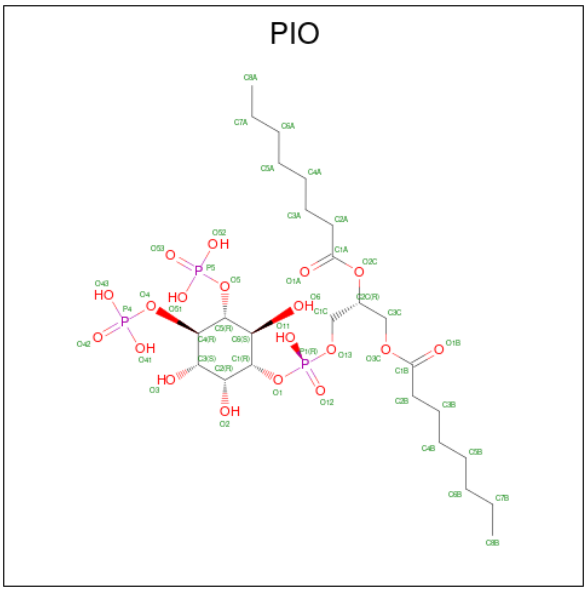


Mol	Chain	Residues	Atoms					AltConf	Trace
8	G	3	Total	C	H	N	O	0	0
			73	22	34	2	15		
8	H	3	Total	C	H	N	O	0	0
			73	22	34	2	15		
8	M	3	Total	C	H	N	O	0	0
			73	22	34	2	15		
8	N	3	Total	C	H	N	O	0	0
			73	22	34	2	15		

- Molecule 9 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.

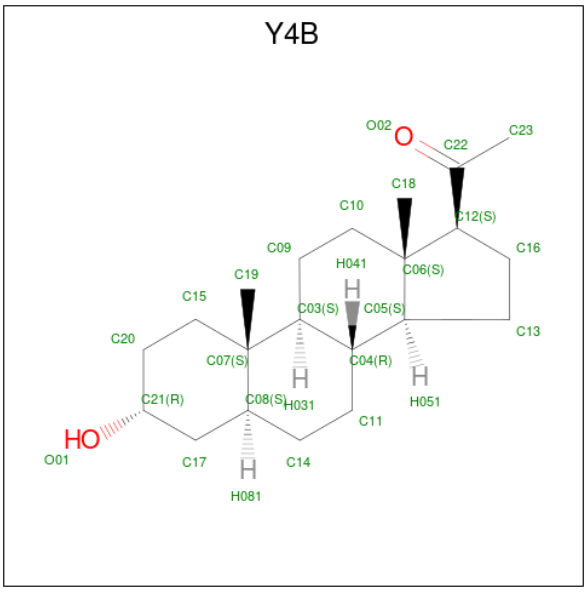
Mol	Chain	Residues	Atoms					AltConf	Trace
9	I	2	Total	C	H	N	O	0	0
			53	16	25	2	10		
9	L	2	Total	C	H	N	O	0	0
			53	16	25	2	10		

- Molecule 10 is [(2R)-2-octanoyloxy-3-[oxidanyl-[(1R,2R,3S,4R,5R,6S)-2,3,6-tris(oxidanyl)-4,5-diphosphonooxy-cyclohexyl]oxy-phosphoryl]oxy-propyl] octanoate (three-letter code: PIO) (formula: $C_{25}H_{49}O_{19}P_3$).



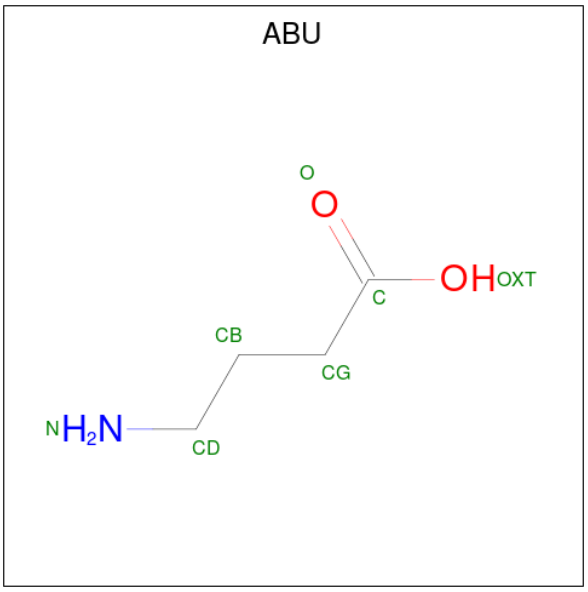
Mol	Chain	Residues	Atoms					AltConf
10	A	1	Total	C	H	O	P	0
			91	25	44	19	3	
10	C	1	Total	C	H	O	P	0
			91	25	44	19	3	

- Molecule 11 is allopregnanolone (three-letter code: Y4B) (formula: $C_{21}H_{34}O_2$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				AltConf
11	A	1	Total	C	H	O	0
			57	21	34	2	
11	B	1	Total	C	H	O	0
			57	21	34	2	

- Molecule 12 is GAMMA-AMINO-BUTANOIC ACID (three-letter code: ABU) (formula: $C_4H_9NO_2$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					AltConf
12	B	1	Total	C	H	N	O	0
			15	4	8	1	2	

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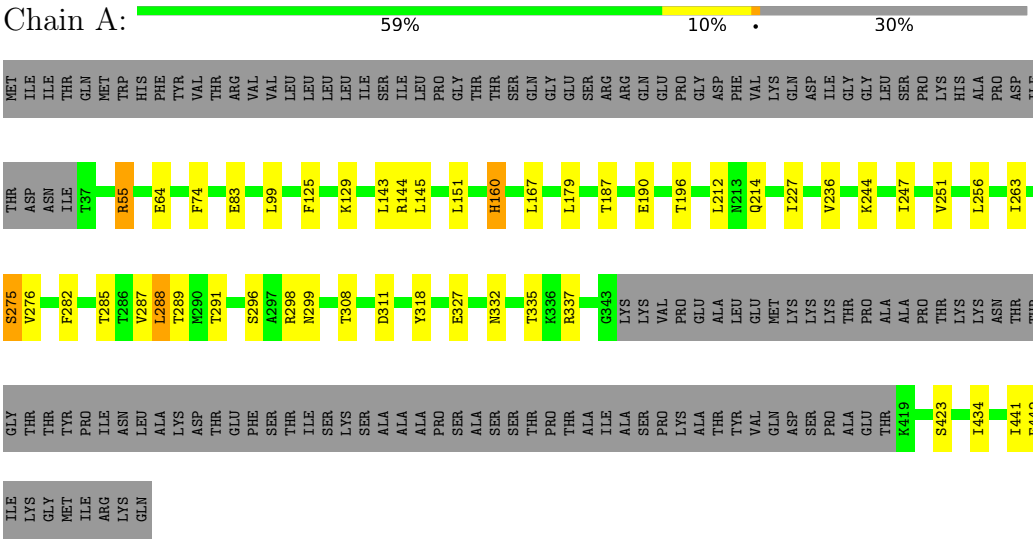
Mol	Chain	Residues	Atoms					AltConf
12	E	1	Total	C	H	N	O	0
			15	4	8	1	2	

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.

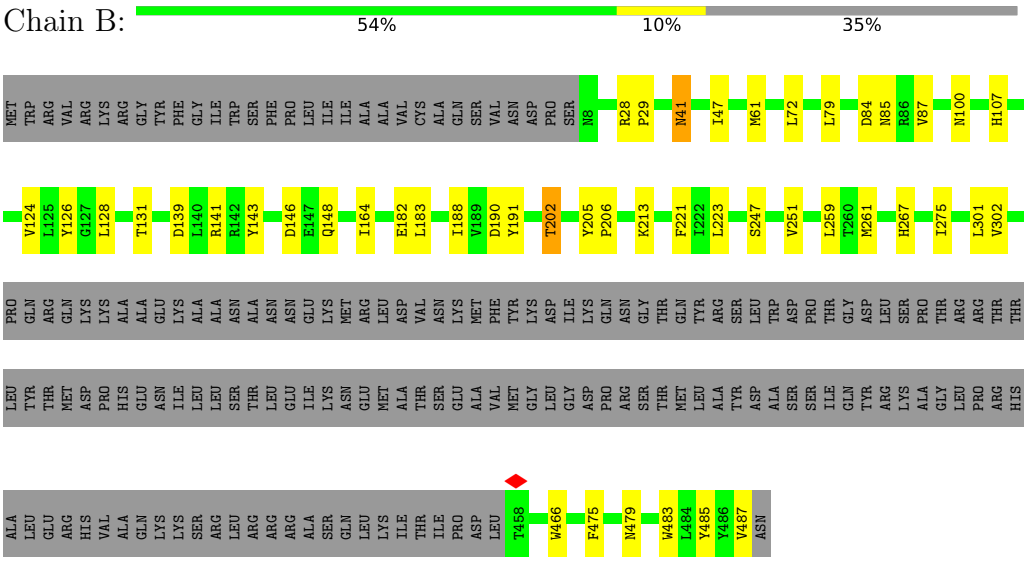
• Molecule 1: Gamma-aminobutyric acid receptor subunit alpha-3

Chain A:



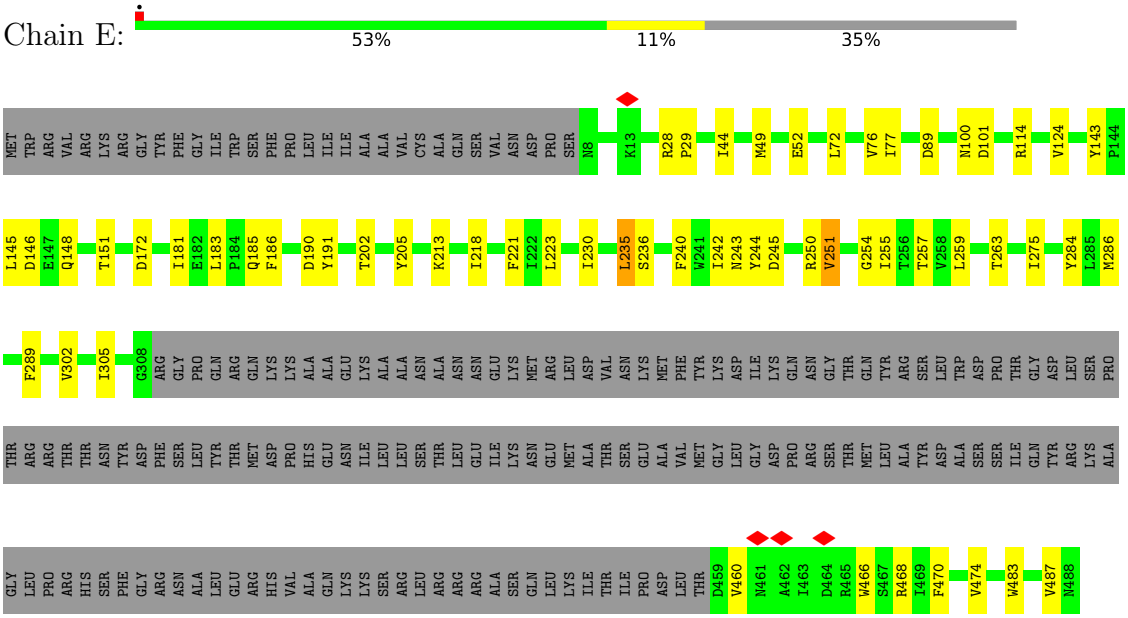
• Molecule 2: Gamma-aminobutyric acid receptor subunit beta-2

Chain B:



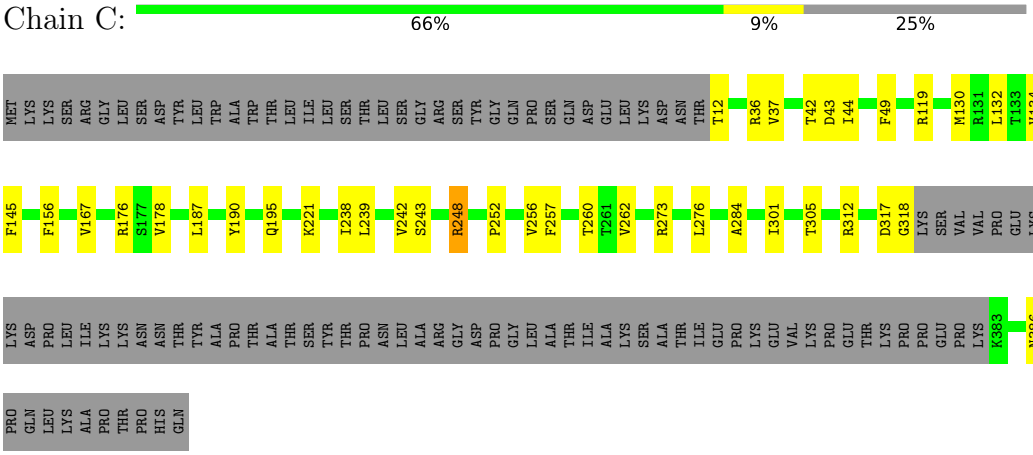
• Molecule 2: Gamma-aminobutyric acid receptor subunit beta-2

Chain E:



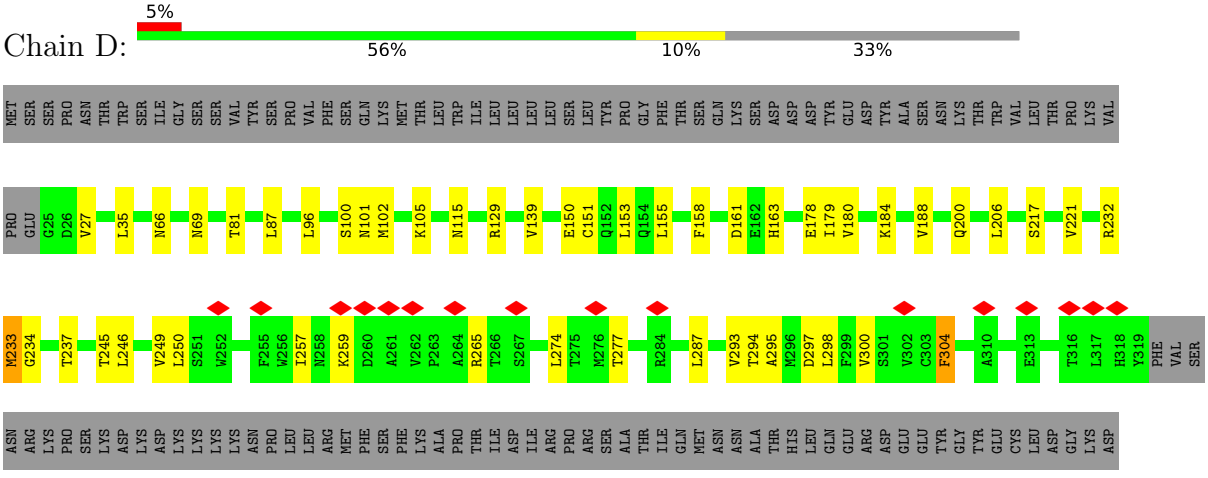
• Molecule 3: Gamma-aminobutyric acid receptor subunit alpha-1

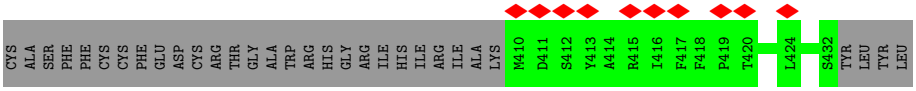
Chain C:



• Molecule 4: Gamma-aminobutyric acid receptor subunit gamma-2

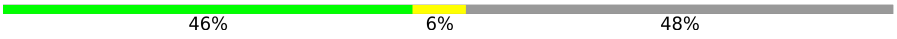
Chain D:

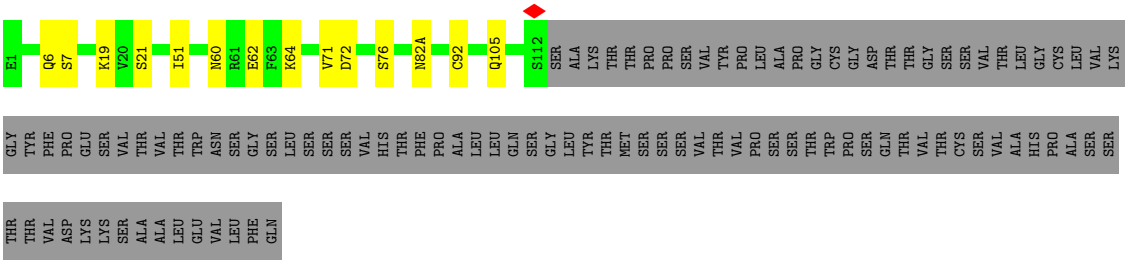




• Molecule 5: Heavy Chain of 8E3 Fab

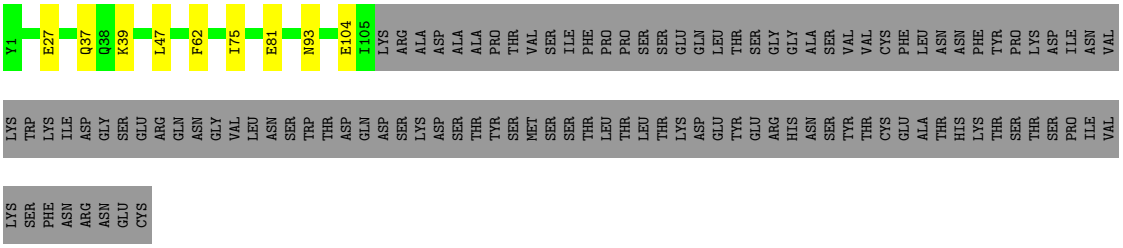
Chain J:





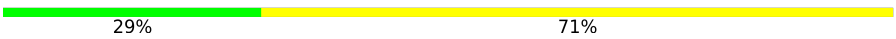
• Molecule 6: Light Chain of 8E3 Fab

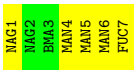
Chain K:



• Molecule 7: alpha-D-mannopyranose-(1-3)-alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose

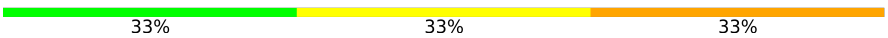
Chain F:





• Molecule 8: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain G:



• Molecule 8: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain H:

• Molecule 8: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain M:

• Molecule 8: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain N:

• Molecule 9: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain I:

• Molecule 9: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain L:

4 Experimental information (i)

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	137566	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/Å^2$)	50	Depositor
Minimum defocus (nm)	800	Depositor
Maximum defocus (nm)	2100	Depositor
Magnification	Not provided	
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	0.730	Depositor
Minimum map value	-0.303	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.019	Depositor
Recommended contour level	0.09	Depositor
Map size (Å)	299.16, 299.16, 299.16	wwPDB
Map dimensions	360, 360, 360	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.83100003, 0.83100003, 0.83100003	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ABU, MAN, FUC, BMA, PIO, NAG, Y4B

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	0.26	0/2850	0.55	0/3879
2	B	0.27	0/2782	0.55	0/3787
2	E	0.27	0/2784	0.55	0/3788
3	C	0.26	0/2867	0.54	0/3895
4	D	0.27	0/2676	0.57	0/3646
5	J	0.28	0/946	0.50	0/1284
6	K	0.29	0/815	0.55	0/1108
All	All	0.27	0/15720	0.55	0/21387

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	2776	2768	2769	28	0
2	B	2712	2717	2716	28	0
2	E	2714	2716	2715	31	0
3	C	2791	2774	2775	24	0
4	D	2605	2582	2582	26	0
5	J	920	863	863	7	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	K	796	756	756	5	0
7	F	82	70	70	0	0
8	G	39	34	34	2	0
8	H	39	34	34	2	0
8	M	39	34	34	0	0
8	N	39	34	34	3	0
9	I	28	25	25	0	0
9	L	28	25	25	0	0
10	A	47	44	44	3	0
10	C	47	44	44	0	0
11	A	23	34	0	0	0
11	B	23	34	0	1	0
12	B	7	8	0	1	0
12	E	7	8	0	0	0
All	All	15762	15604	15520	147	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 5.

All (147) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:296:SER:O	1:A:299:ASN:ND2	2.19	0.76
2:B:213:LYS:HB2	8:H:1:NAG:H81	1.68	0.75
2:B:259:LEU:HD11	3:C:260:THR:HG23	1.69	0.74
2:E:146:ASP:OD2	2:E:148:GLN:NE2	2.21	0.73
2:E:302:VAL:HG13	2:E:460:VAL:HG13	1.71	0.71
2:E:185:GLN:N	2:E:185:GLN:OE1	2.23	0.71
2:E:259:LEU:O	2:E:263:THR:HG23	1.91	0.70
2:B:182:GLU:N	2:B:182:GLU:OE1	2.23	0.70
4:D:101:ASN:O	4:D:105:LYS:NZ	2.22	0.69
4:D:35:LEU:HD21	4:D:102:MET:HE3	1.74	0.68
2:B:483:TRP:O	2:B:487:VAL:N	2.28	0.67
5:J:62:GLU:OE1	5:J:62:GLU:N	2.28	0.66
2:E:470:PHE:O	2:E:474:VAL:HG12	1.97	0.65
2:B:28:ARG:NH1	2:B:29:PRO:O	2.30	0.65
4:D:188:VAL:HB	4:D:206:LEU:HD21	1.78	0.64
3:C:144:ASP:OD2	3:C:417:ARG:NH1	2.30	0.63
1:A:227:ILE:HD12	1:A:236:VAL:HG21	1.81	0.63
1:A:144:ARG:NH2	2:E:205:TYR:OH	2.32	0.62
3:C:43:ASP:OD1	3:C:44:ILE:N	2.33	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:66:ASN:ND2	4:D:69:ASN:OD1	2.33	0.61
2:E:190:ASP:OD1	2:E:191:TYR:N	2.33	0.61
2:B:41:ASN:N	2:B:41:ASN:OD1	2.32	0.60
3:C:248:ARG:NH1	3:C:387:SER:O	2.35	0.60
2:B:146:ASP:OD2	2:B:148:GLN:NE2	2.35	0.59
2:E:483:TRP:O	2:E:487:VAL:N	2.35	0.59
2:E:100:ASN:ND2	2:E:151:THR:O	2.35	0.59
1:A:332:ASN:O	1:A:335:THR:OG1	2.20	0.59
1:A:167:LEU:O	1:A:308:THR:HG22	2.04	0.58
1:A:287:VAL:O	1:A:291:THR:HG23	2.03	0.58
1:A:288:LEU:O	1:A:291:THR:OG1	2.20	0.58
8:H:2:NAG:H83	8:H:2:NAG:H3	1.85	0.58
2:B:259:LEU:HD11	3:C:260:THR:CG2	2.33	0.57
2:E:230:ILE:HD11	2:E:284:TYR:OH	2.03	0.57
8:G:2:NAG:H83	8:G:2:NAG:H3	1.87	0.56
2:E:44:ILE:HG22	2:E:181:ILE:HD11	1.87	0.56
4:D:178:GLU:O	4:D:179:ILE:HD12	2.05	0.56
2:B:475:PHE:O	2:B:479:ASN:ND2	2.39	0.55
1:A:99:LEU:HD12	1:A:151:LEU:HD11	1.89	0.55
1:A:263:ILE:O	1:A:267:VAL:HG23	2.07	0.55
1:A:337:ARG:NH1	10:A:501:PIO:O52	2.40	0.54
3:C:130:MET:SD	3:C:132:LEU:HD21	2.47	0.54
6:K:104:GLU:N	6:K:104:GLU:OE1	2.41	0.54
3:C:262:VAL:HG11	4:D:250:LEU:HD11	1.89	0.54
2:E:52:GLU:N	2:E:52:GLU:OE1	2.38	0.54
2:E:76:VAL:HG23	2:E:77:ILE:HG23	1.90	0.54
3:C:176:ARG:NH1	6:K:93:ASN:OD1	2.41	0.53
4:D:234:GLY:O	4:D:237:THR:OG1	2.23	0.53
1:A:273:ARG:NH1	10:A:501:PIO:O6	2.41	0.53
3:C:190:TYR:O	3:C:221:LYS:NZ	2.40	0.53
2:B:47:ILE:HD12	2:B:188:ILE:HD11	1.90	0.53
2:B:126:TYR:OH	2:B:128:LEU:HD21	2.09	0.53
4:D:151:CYS:O	4:D:153:LEU:HD12	2.08	0.53
2:B:139:ASP:OD1	2:B:141:ARG:NE	2.42	0.52
2:B:190:ASP:OD1	2:B:191:TYR:N	2.43	0.52
8:N:2:NAG:H3	8:N:2:NAG:H83	1.90	0.52
1:A:291:THR:HG22	1:A:318:TYR:OH	2.09	0.52
2:E:240:PHE:O	2:E:468:ARG:NH1	2.42	0.52
10:A:501:PIO:P1	10:A:501:PIO:HO2	2.33	0.52
2:B:72:LEU:HD12	2:B:124:VAL:HG21	1.91	0.51
6:K:62:PHE:CD1	6:K:75:ILE:HD12	2.45	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:55:ARG:NE	1:A:190:GLU:O	2.44	0.51
4:D:158:PHE:O	4:D:295:ALA:HB3	2.10	0.51
5:J:6:GLN:OE1	5:J:92:CYS:N	2.44	0.50
4:D:180:VAL:HA	4:D:221:VAL:HG21	1.92	0.50
1:A:74:PHE:CD1	1:A:212:LEU:HD21	2.46	0.50
4:D:27:VAL:HG13	4:D:96:LEU:HD22	1.94	0.50
2:E:183:LEU:HD13	2:E:186:PHE:HB2	1.93	0.50
2:E:213:LYS:HD3	8:N:1:NAG:H82	1.93	0.50
6:K:39:LYS:NZ	6:K:81:GLU:OE2	2.43	0.50
1:A:125:PHE:HE2	1:A:179:LEU:HD11	1.77	0.49
1:A:144:ARG:NH1	2:E:202:THR:O	2.43	0.49
2:E:243:ASN:ND2	2:E:245:ASP:OD1	2.46	0.49
2:E:254:GLY:O	2:E:257:THR:OG1	2.27	0.48
2:B:275:ILE:HD12	2:B:275:ILE:O	2.14	0.48
3:C:42:THR:HG21	3:C:156:PHE:CZ	2.49	0.48
2:B:205:TYR:OH	3:C:119:ARG:NH2	2.45	0.48
3:C:301:ILE:O	3:C:305:THR:HG23	2.14	0.48
2:B:107:HIS:NE2	2:B:131:THR:OG1	2.38	0.48
2:E:275:ILE:HD12	2:E:275:ILE:O	2.14	0.48
3:C:239:LEU:O	3:C:242:VAL:HG22	2.14	0.47
5:J:72:ASP:O	5:J:76:SER:N	2.47	0.47
4:D:245:THR:O	4:D:249:VAL:HG23	2.13	0.47
2:E:242:ILE:O	2:E:250:ARG:NH1	2.47	0.47
1:A:273:ARG:NH1	1:A:423:SER:O	2.46	0.47
2:E:183:LEU:HD13	2:E:186:PHE:CB	2.44	0.47
4:D:100:SER:HG	4:D:129:ARG:HE	1.61	0.47
1:A:275:SER:O	1:A:275:SER:OG	2.30	0.47
4:D:87:LEU:HD12	4:D:139:VAL:HG21	1.97	0.47
2:B:84:ASP:OD1	2:B:85:ASN:N	2.48	0.46
4:D:233:MET:O	4:D:237:THR:HG23	2.15	0.46
5:J:60:ASN:O	5:J:64:LYS:N	2.48	0.46
3:C:238:ILE:O	3:C:242:VAL:HG13	2.16	0.46
4:D:115:ASN:O	4:D:150:GLU:N	2.43	0.46
3:C:49:PHE:CD2	3:C:187:LEU:HD21	2.50	0.46
3:C:317:ASP:OD1	3:C:318:GLY:N	2.48	0.46
6:K:37:GLN:HB2	6:K:47:LEU:HD11	1.98	0.46
4:D:161:ASP:OD2	4:D:163:HIS:NE2	2.46	0.45
2:B:301:LEU:HD23	11:B:502:Y4B:O02	2.15	0.45
3:C:312:ARG:O	3:C:386:ASN:ND2	2.48	0.45
4:D:155:LEU:O	4:D:294:THR:HG22	2.16	0.45
2:B:85:ASN:O	2:B:114:ARG:NH2	2.47	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:178:VAL:O	3:C:195:GLN:NE2	2.50	0.45
2:E:236:SER:CB	2:E:257:THR:HG21	2.46	0.45
5:J:7:SER:OG	5:J:21:SER:OG	2.35	0.45
1:A:129:LYS:N	1:A:160:HIS:O	2.47	0.45
1:A:83:GLU:OE1	1:A:160:HIS:ND1	2.50	0.45
2:B:202:THR:HG22	3:C:119:ARG:CZ	2.46	0.44
2:B:84:ASP:O	2:B:87:VAL:HG12	2.18	0.44
3:C:252:PRO:O	3:C:256:VAL:HG23	2.17	0.44
2:E:143:TYR:HH	2:E:221:PHE:HE1	1.65	0.44
8:N:1:NAG:O7	8:N:1:NAG:O3	2.35	0.44
1:A:274:GLU:OE1	1:A:274:GLU:N	2.49	0.44
4:D:200:GLN:O	4:D:232:ARG:N	2.45	0.44
2:E:28:ARG:NH1	2:E:29:PRO:O	2.51	0.44
8:G:2:NAG:O3	8:G:3:BMA:O5	2.31	0.44
3:C:142:LEU:HD13	3:C:276:LEU:HD22	2.00	0.44
3:C:145:PHE:O	3:C:284:ALA:HB3	2.18	0.43
2:B:247:SER:O	2:B:251:VAL:HG22	2.18	0.43
3:C:49:PHE:HD2	3:C:187:LEU:HD21	1.83	0.43
3:C:37:VAL:HG11	3:C:167:VAL:HG23	2.00	0.43
2:E:89:ASP:OD1	2:E:114:ARG:NH1	2.52	0.43
5:J:6:GLN:O	5:J:105:GLN:NE2	2.52	0.43
4:D:293:VAL:HG13	4:D:293:VAL:O	2.19	0.43
4:D:297:ASP:HA	4:D:300:VAL:HG22	2.00	0.43
1:A:143:LEU:HD21	1:A:151:LEU:HD23	2.01	0.42
1:A:64:GLU:OE1	1:A:64:GLU:N	2.53	0.42
2:B:164:ILE:HG23	2:B:206:PRO:HG3	2.01	0.42
4:D:277:THR:HG22	4:D:304:PHE:CE2	2.55	0.42
4:D:274:LEU:O	4:D:277:THR:OG1	2.28	0.41
2:E:145:LEU:HD11	2:E:218:ILE:HD12	2.01	0.41
1:A:441:ILE:HD12	1:A:442:PHE:N	2.35	0.41
5:J:51:ILE:HD13	5:J:71:VAL:HG13	2.03	0.41
2:E:251:VAL:HG13	2:E:255:ILE:HD12	2.02	0.41
1:A:285:THR:O	1:A:289:THR:OG1	2.29	0.41
1:A:298:ARG:NH1	1:A:311:ASP:OD2	2.53	0.41
4:D:257:ILE:O	4:D:265:ARG:NH2	2.44	0.41
2:E:145:LEU:CD1	2:E:218:ILE:HD12	2.50	0.41
1:A:247:ILE:O	1:A:251:VAL:HG23	2.20	0.41
1:A:99:LEU:HB3	1:A:145:LEU:HD11	2.02	0.41
2:B:202:THR:OG1	12:B:501:ABU:O	2.32	0.41
2:B:143:TYR:HH	2:B:221:PHE:HE1	1.68	0.41
2:E:235:LEU:HD12	2:E:235:LEU:C	2.40	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:111:VAL:HG22	2:B:112:LYS:H	1.87	0.41
2:E:72:LEU:HD12	2:E:124:VAL:HG21	2.03	0.40
4:D:246:LEU:O	4:D:250:LEU:HD13	2.21	0.40
2:B:126:TYR:CZ	2:B:128:LEU:HD21	2.56	0.40
4:D:81:THR:HG21	4:D:184:LYS:NZ	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	339/492 (69%)	330 (97%)	9 (3%)	0	100	100
2	B	327/512 (64%)	317 (97%)	10 (3%)	0	100	100
2	E	327/512 (64%)	317 (97%)	10 (3%)	0	100	100
3	C	340/455 (75%)	331 (97%)	9 (3%)	0	100	100
4	D	314/474 (66%)	305 (97%)	9 (3%)	0	100	100
5	J	115/223 (52%)	111 (96%)	4 (4%)	0	100	100
6	K	103/213 (48%)	99 (96%)	4 (4%)	0	100	100
All	All	1865/2881 (65%)	1810 (97%)	55 (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	307/435 (71%)	294 (96%)	13 (4%)	30	45
2	B	301/457 (66%)	287 (95%)	14 (5%)	26	40
2	E	301/457 (66%)	290 (96%)	11 (4%)	34	51
3	C	306/404 (76%)	299 (98%)	7 (2%)	50	68
4	D	292/436 (67%)	286 (98%)	6 (2%)	53	71
5	J	98/195 (50%)	96 (98%)	2 (2%)	55	72
6	K	83/188 (44%)	82 (99%)	1 (1%)	71	83
All	All	1688/2572 (66%)	1634 (97%)	54 (3%)	42	56

All (54) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	55	ARG
1	A	160	HIS
1	A	187	THR
1	A	196	THR
1	A	214	GLN
1	A	244	LYS
1	A	256	LEU
1	A	275	SER
1	A	276	VAL
1	A	282	PHE
1	A	288	LEU
1	A	327	GLU
1	A	434	ILE
2	B	41	ASN
2	B	61	MET
2	B	79	LEU
2	B	100	ASN
2	B	115	MET
2	B	183	LEU
2	B	202	THR
2	B	223	LEU
2	B	261	MET
2	B	267	HIS
2	B	302	VAL
2	B	305	ILE
2	B	466	TRP
2	B	485	TYR
3	C	12	THR
3	C	36	ARG

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Mol	Chain	Res	Type
3	C	134	VAL
3	C	243	SER
3	C	248	ARG
3	C	257	PHE
3	C	273	ARG
4	D	217	SER
4	D	233	MET
4	D	259	LYS
4	D	287	LEU
4	D	298	LEU
4	D	304	PHE
2	E	49	MET
2	E	101	ASP
2	E	172	ASP
2	E	223	LEU
2	E	235	LEU
2	E	244	TYR
2	E	251	VAL
2	E	286	MET
2	E	289	PHE
2	E	305	ILE
2	E	466	TRP
5	J	19	LYS
5	J	82(A)	ASN
6	K	27	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	175	HIS
1	A	214	GLN
1	A	299	ASN
1	A	443	ASN
2	B	64	GLN
2	B	85	ASN
2	B	185	GLN
3	C	247	ASN
4	D	66	ASN
4	D	69	ASN
5	J	82(A)	ASN
5	J	98	ASN
6	K	79	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

23 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	$\#\|Z\| > 2$
7	NAG	F	1	7,1	14,14,15	1.14	1 (7%)	17,19,21	2.17	1 (5%)
7	NAG	F	2	7	14,14,15	0.22	0	17,19,21	0.64	0
7	BMA	F	3	7	11,11,12	0.54	0	15,15,17	0.80	0
7	MAN	F	4	7	11,11,12	0.84	1 (9%)	15,15,17	1.27	2 (13%)
7	MAN	F	5	7	11,11,12	0.69	0	15,15,17	1.39	2 (13%)
7	MAN	F	6	7	11,11,12	0.57	0	15,15,17	1.02	2 (13%)
7	FUC	F	7	7	10,10,11	0.74	0	14,14,16	0.90	1 (7%)
8	NAG	G	1	8,2	14,14,15	0.39	0	17,19,21	0.71	0
8	NAG	G	2	8	14,14,15	0.28	0	17,19,21	1.17	1 (5%)
8	BMA	G	3	8	11,11,12	0.52	0	15,15,17	0.71	0
8	NAG	H	1	8,2	14,14,15	0.31	0	17,19,21	0.75	1 (5%)
8	NAG	H	2	8	14,14,15	0.27	0	17,19,21	1.18	1 (5%)
8	BMA	H	3	8	11,11,12	0.96	1 (9%)	15,15,17	0.86	1 (6%)
9	NAG	I	1	9,3	14,14,15	0.14	0	17,19,21	0.71	1 (5%)
9	NAG	I	2	9	14,14,15	0.17	0	17,19,21	0.47	0
9	NAG	L	1	9,4	14,14,15	0.58	0	17,19,21	1.44	3 (17%)
9	NAG	L	2	9	14,14,15	0.22	0	17,19,21	0.42	0
8	NAG	M	1	8,2	14,14,15	0.41	0	17,19,21	0.67	1 (5%)
8	NAG	M	2	8	14,14,15	0.22	0	17,19,21	0.40	0
8	BMA	M	3	8	11,11,12	0.52	0	15,15,17	0.75	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
8	NAG	N	1	8,2	14,14,15	0.53	0	17,19,21	0.69	1 (5%)
8	NAG	N	2	8	14,14,15	0.26	0	17,19,21	1.19	1 (5%)
8	BMA	N	3	8	11,11,12	0.54	0	15,15,17	0.83	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
7	NAG	F	1	7,1	-	0/6/23/26	0/1/1/1
7	NAG	F	2	7	-	4/6/23/26	0/1/1/1
7	BMA	F	3	7	-	2/2/19/22	0/1/1/1
7	MAN	F	4	7	-	0/2/19/22	1/1/1/1
7	MAN	F	5	7	-	2/2/19/22	1/1/1/1
7	MAN	F	6	7	-	1/2/19/22	0/1/1/1
7	FUC	F	7	7	-	-	0/1/1/1
8	NAG	G	1	8,2	-	4/6/23/26	0/1/1/1
8	NAG	G	2	8	-	5/6/23/26	0/1/1/1
8	BMA	G	3	8	-	0/2/19/22	0/1/1/1
8	NAG	H	1	8,2	-	3/6/23/26	0/1/1/1
8	NAG	H	2	8	-	5/6/23/26	0/1/1/1
8	BMA	H	3	8	-	2/2/19/22	0/1/1/1
9	NAG	I	1	9,3	-	4/6/23/26	0/1/1/1
9	NAG	I	2	9	-	4/6/23/26	0/1/1/1
9	NAG	L	1	9,4	-	4/6/23/26	0/1/1/1
9	NAG	L	2	9	-	0/6/23/26	0/1/1/1
8	NAG	M	1	8,2	-	2/6/23/26	0/1/1/1
8	NAG	M	2	8	-	1/6/23/26	0/1/1/1
8	BMA	M	3	8	-	1/2/19/22	0/1/1/1
8	NAG	N	1	8,2	-	4/6/23/26	0/1/1/1
8	NAG	N	2	8	-	4/6/23/26	0/1/1/1
8	BMA	N	3	8	-	0/2/19/22	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	F	1	NAG	O5-C1	4.17	1.50	1.43
8	H	3	BMA	C1-C2	2.87	1.58	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	F	4	MAN	C1-C2	2.30	1.57	1.52

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
7	F	1	NAG	C1-O5-C5	8.60	123.85	112.19
9	L	1	NAG	C1-O5-C5	4.48	118.26	112.19
7	F	5	MAN	C1-O5-C5	4.20	117.89	112.19
8	N	2	NAG	C2-N2-C7	3.85	128.39	122.90
8	G	2	NAG	C2-N2-C7	3.85	128.38	122.90
8	H	2	NAG	C2-N2-C7	3.84	128.37	122.90
7	F	4	MAN	C1-O5-C5	3.71	117.22	112.19
9	I	1	NAG	C1-O5-C5	2.59	115.70	112.19
7	F	6	MAN	C1-O5-C5	2.51	115.59	112.19
8	N	1	NAG	C1-O5-C5	2.47	115.54	112.19
7	F	4	MAN	O2-C2-C3	-2.47	105.20	110.14
7	F	5	MAN	O2-C2-C3	-2.31	105.50	110.14
8	H	1	NAG	C1-O5-C5	2.27	115.27	112.19
9	L	1	NAG	C3-C4-C5	2.24	114.24	110.24
8	M	1	NAG	C2-N2-C7	2.21	126.05	122.90
7	F	6	MAN	O2-C2-C3	-2.16	105.81	110.14
9	L	1	NAG	C2-N2-C7	2.11	125.91	122.90
8	H	3	BMA	O2-C2-C3	-2.09	105.95	110.14
7	F	7	FUC	C1-C2-C3	2.05	112.18	109.67

There are no chirality outliers.

All (52) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	F	5	MAN	O5-C5-C6-O6
8	H	3	BMA	O5-C5-C6-O6
9	L	1	NAG	O5-C5-C6-O6
7	F	5	MAN	C4-C5-C6-O6
9	I	1	NAG	O5-C5-C6-O6
8	H	3	BMA	C4-C5-C6-O6
8	H	1	NAG	O5-C5-C6-O6
7	F	2	NAG	O5-C5-C6-O6
8	G	1	NAG	O5-C5-C6-O6
7	F	3	BMA	C4-C5-C6-O6
9	L	1	NAG	C4-C5-C6-O6
7	F	2	NAG	C8-C7-N2-C2
7	F	2	NAG	O7-C7-N2-C2

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Mol	Chain	Res	Type	Atoms
8	G	1	NAG	C8-C7-N2-C2
8	G	1	NAG	O7-C7-N2-C2
8	G	2	NAG	C8-C7-N2-C2
8	G	2	NAG	O7-C7-N2-C2
8	H	2	NAG	C8-C7-N2-C2
8	H	2	NAG	O7-C7-N2-C2
8	M	1	NAG	C8-C7-N2-C2
8	M	1	NAG	O7-C7-N2-C2
8	N	2	NAG	C8-C7-N2-C2
8	N	2	NAG	O7-C7-N2-C2
9	L	1	NAG	C8-C7-N2-C2
9	L	1	NAG	O7-C7-N2-C2
8	H	1	NAG	C4-C5-C6-O6
9	I	1	NAG	C4-C5-C6-O6
8	G	1	NAG	C4-C5-C6-O6
7	F	2	NAG	C4-C5-C6-O6
8	N	1	NAG	O5-C5-C6-O6
7	F	3	BMA	O5-C5-C6-O6
8	N	1	NAG	C1-C2-N2-C7
8	M	3	BMA	O5-C5-C6-O6
8	H	2	NAG	C4-C5-C6-O6
8	G	2	NAG	C4-C5-C6-O6
8	M	2	NAG	O5-C5-C6-O6
7	F	6	MAN	O5-C5-C6-O6
8	H	2	NAG	O5-C5-C6-O6
8	N	1	NAG	C4-C5-C6-O6
8	G	2	NAG	O5-C5-C6-O6
8	N	1	NAG	C3-C2-N2-C7
8	N	2	NAG	C3-C2-N2-C7
9	I	2	NAG	C4-C5-C6-O6
9	I	2	NAG	C1-C2-N2-C7
9	I	1	NAG	C1-C2-N2-C7
8	G	2	NAG	C3-C2-N2-C7
8	H	1	NAG	C3-C2-N2-C7
8	H	2	NAG	C3-C2-N2-C7
9	I	1	NAG	C3-C2-N2-C7
9	I	2	NAG	C3-C2-N2-C7
9	I	2	NAG	O5-C5-C6-O6
8	N	2	NAG	C1-C2-N2-C7

All (2) ring outliers are listed below:

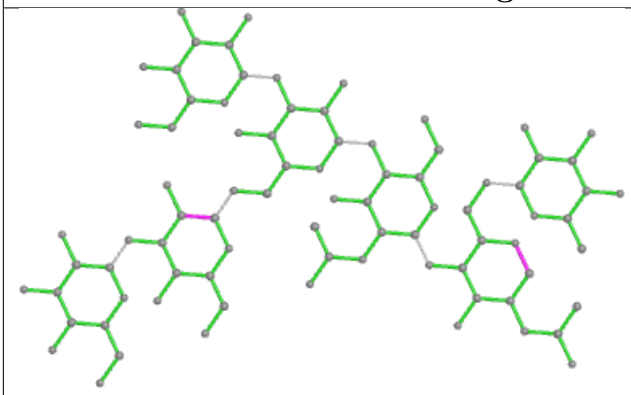
Mol	Chain	Res	Type	Atoms
7	F	5	MAN	C1-C2-C3-C4-C5-O5
7	F	4	MAN	C1-C2-C3-C4-C5-O5

6 monomers are involved in 7 short contacts:

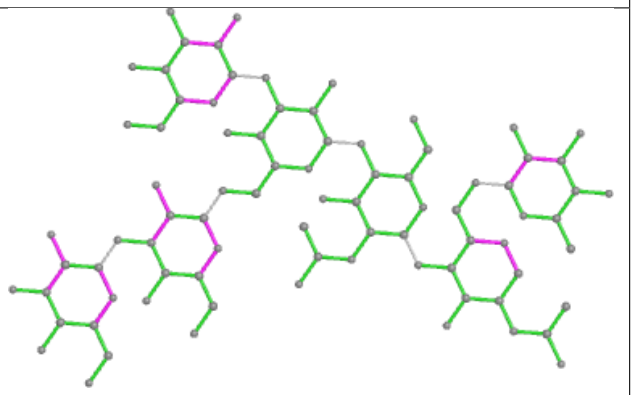
Mol	Chain	Res	Type	Clashes	Symm-Clashes
8	G	3	BMA	1	0
8	G	2	NAG	2	0
8	H	2	NAG	1	0
8	H	1	NAG	1	0
8	N	1	NAG	2	0
8	N	2	NAG	1	0

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

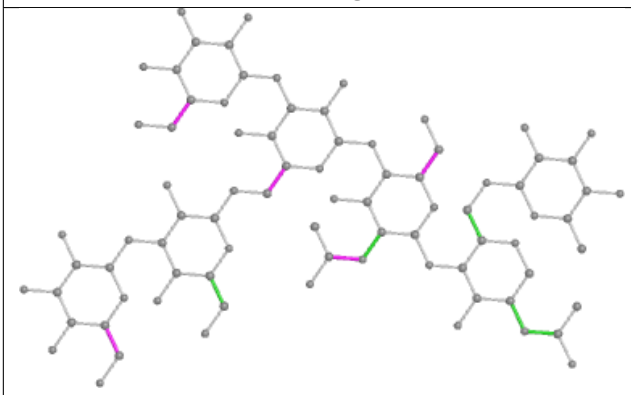
Oligosaccharide Chain F



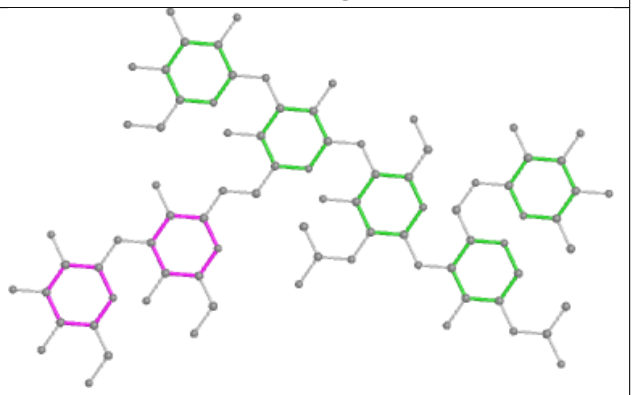
Bond lengths



Bond angles

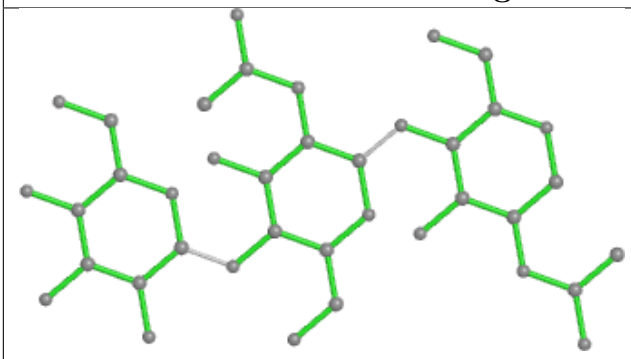


Torsions

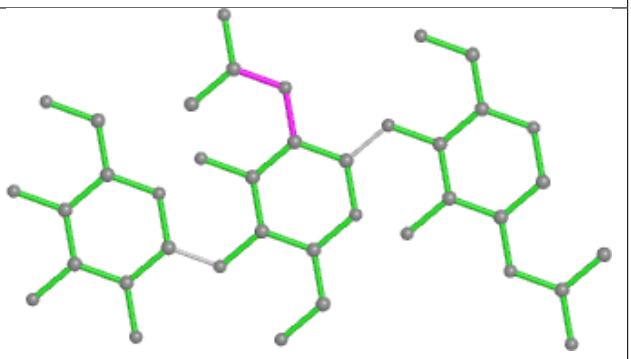


Rings

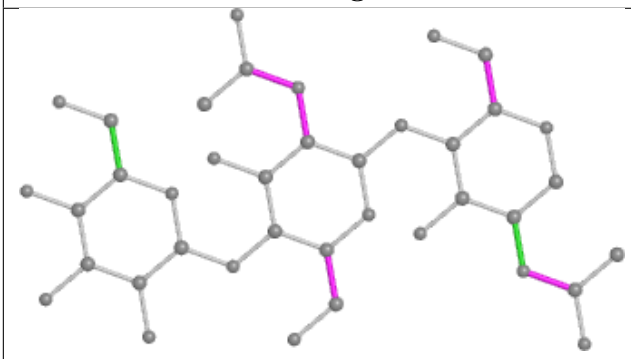
Oligosaccharide Chain G



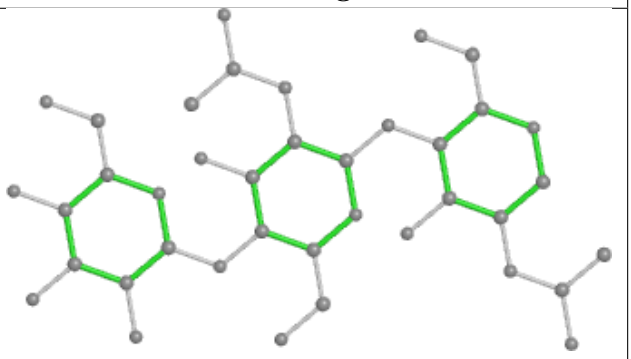
Bond lengths



Bond angles

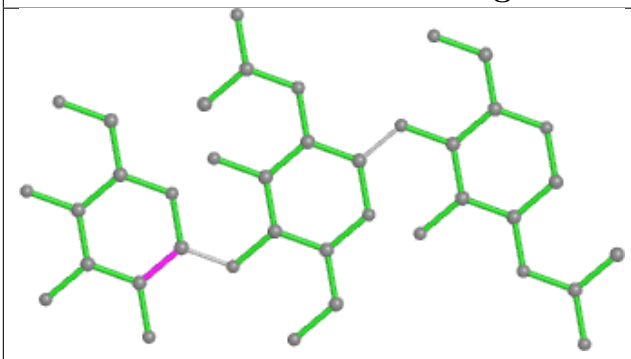


Torsions

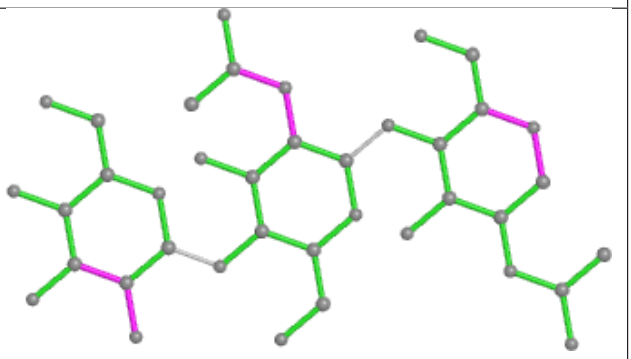


Rings

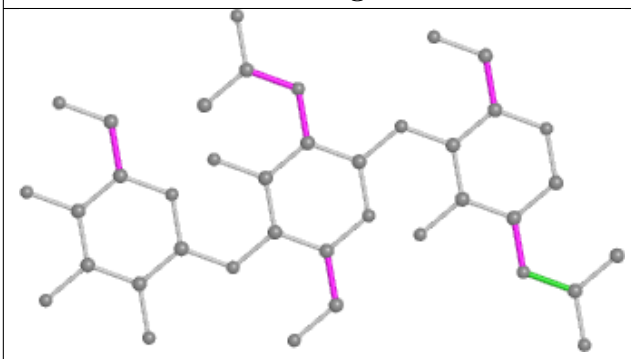
Oligosaccharide Chain H



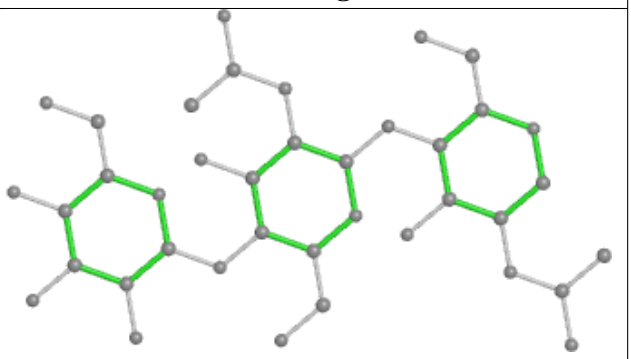
Bond lengths



Bond angles

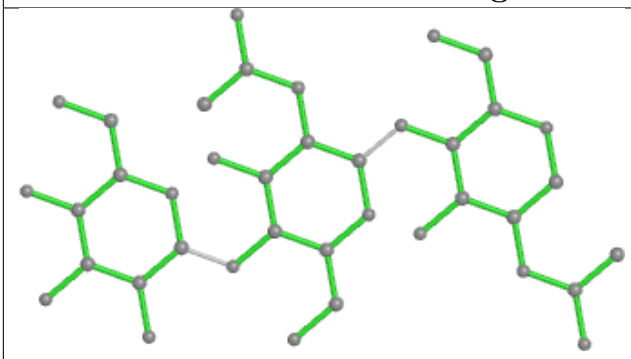


Torsions

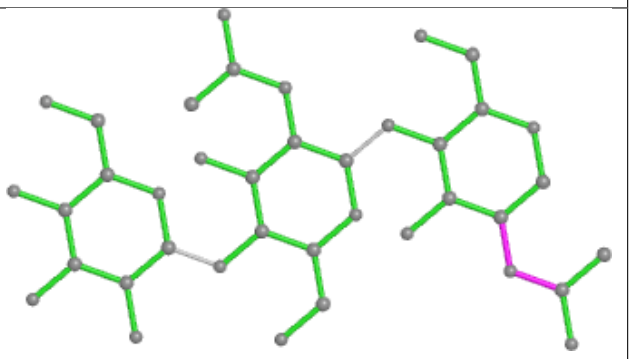


Rings

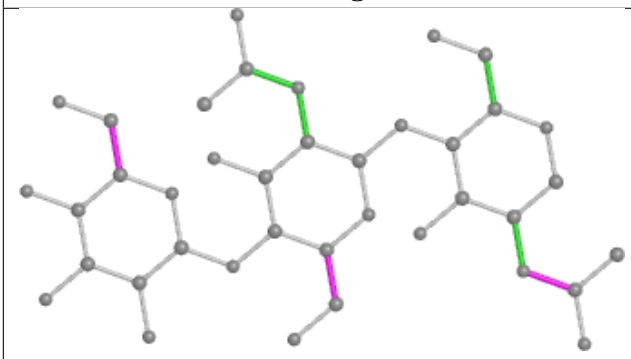
Oligosaccharide Chain M



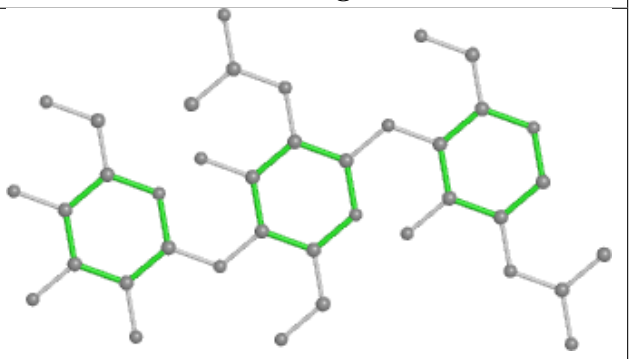
Bond lengths



Bond angles

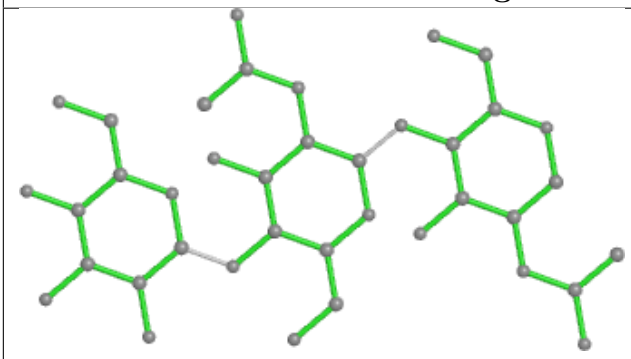


Torsions

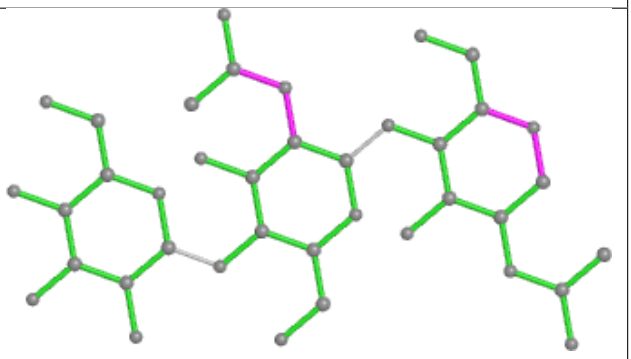


Rings

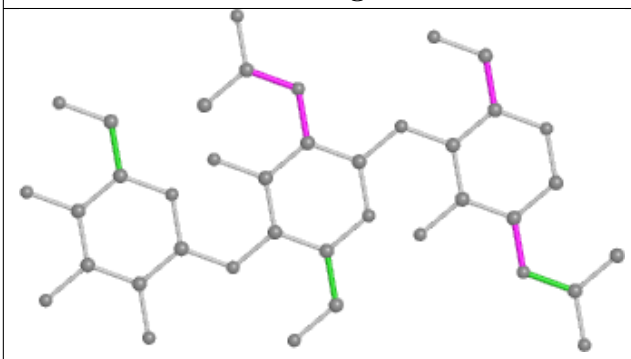
Oligosaccharide Chain N



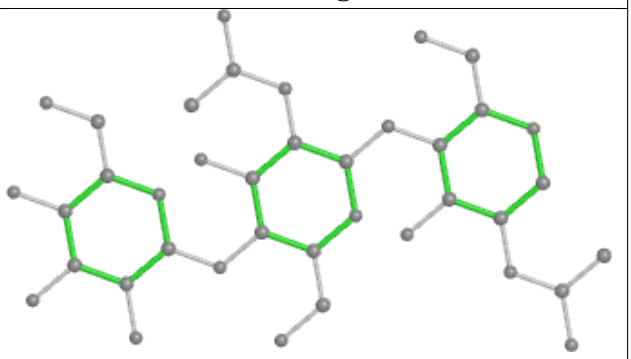
Bond lengths



Bond angles

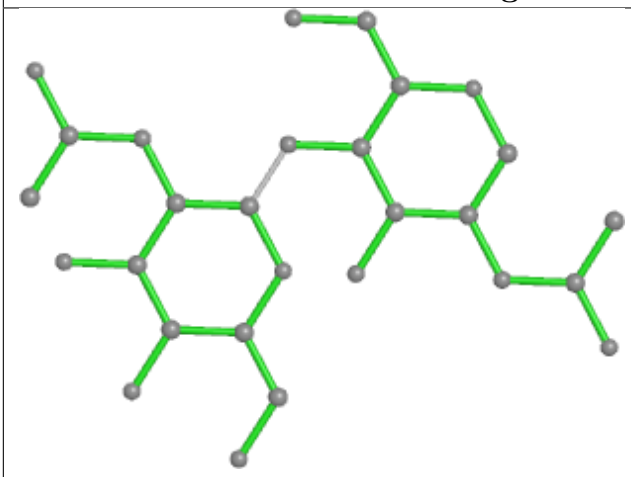


Torsions

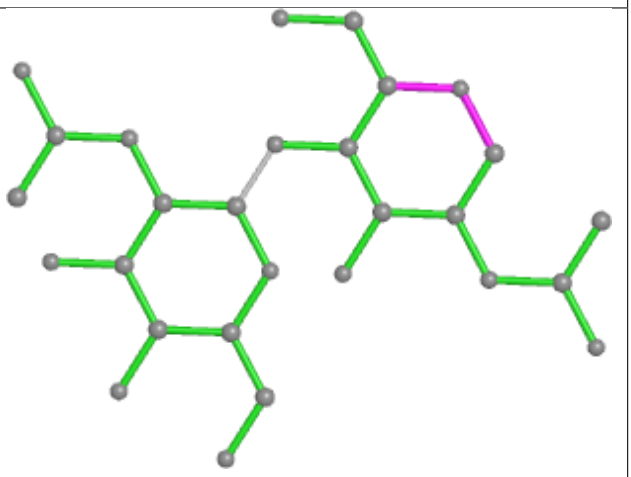


Rings

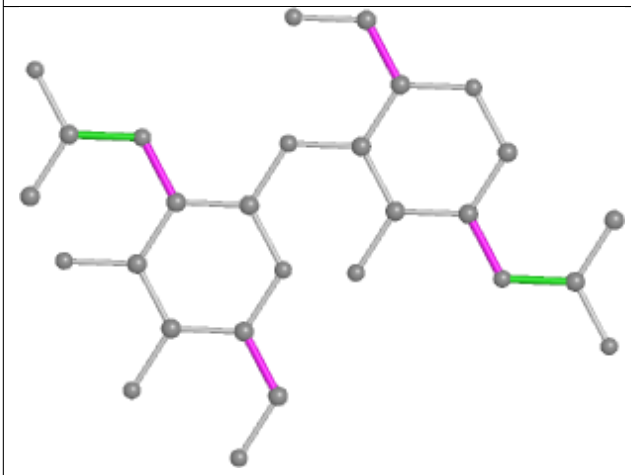
Oligosaccharide Chain I



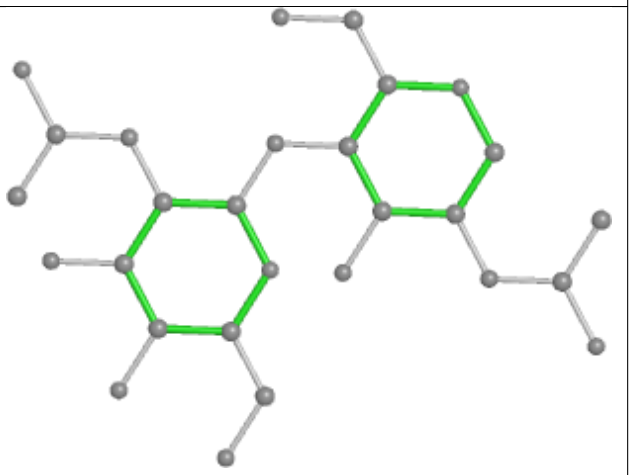
Bond lengths



Bond angles

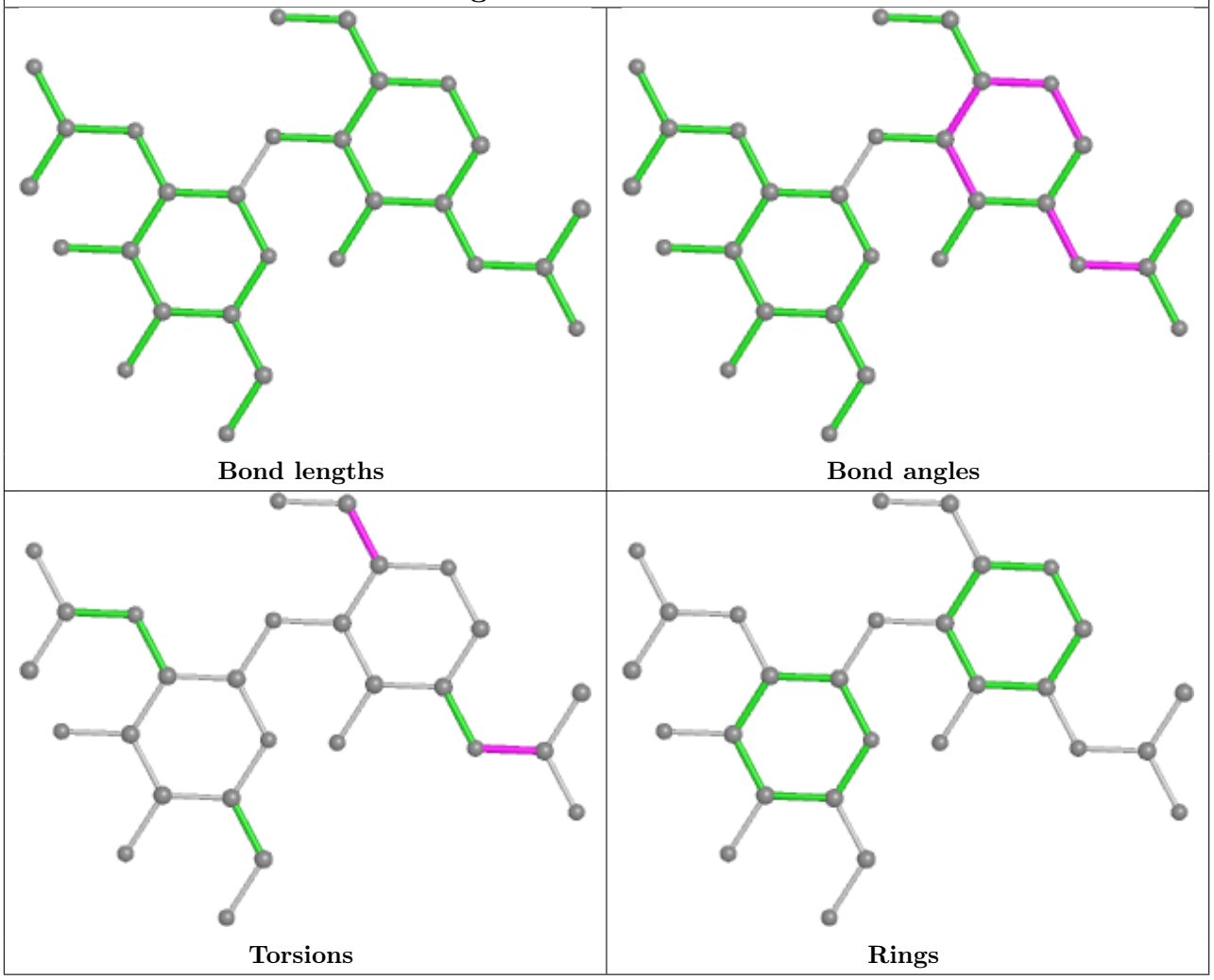


Torsions



Rings

Oligosaccharide Chain L



5.6 Ligand geometry (i)

6 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	$\#\|Z\| > 2$
12	ABU	E	501	-	6,6,6	0.86	0	6,6,6	1.40	1 (16%)
12	ABU	B	501	-	6,6,6	0.87	0	6,6,6	1.25	0
11	Y4B	B	502	-	26,26,26	1.52	3 (11%)	42,42,42	2.19	17 (40%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
10	PIO	C	601	-	47,47,47	1.37	10 (21%)	61,65,65	1.13	3 (4%)
11	Y4B	A	502	-	26,26,26	1.70	4 (15%)	42,42,42	2.26	16 (38%)
10	PIO	A	501	-	47,47,47	1.37	10 (21%)	61,65,65	1.19	4 (6%)

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
12	ABU	E	501	-	-	2/4/4/4	-
12	ABU	B	501	-	-	2/4/4/4	-
11	Y4B	B	502	-	-	0/4/62/62	0/4/4/4
10	PIO	C	601	-	-	16/44/68/68	0/1/1/1
11	Y4B	A	502	-	-	0/4/62/62	0/4/4/4
10	PIO	A	501	-	-	13/44/68/68	0/1/1/1

All (27) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
11	A	502	Y4B	C04-C05	4.98	1.63	1.53
11	A	502	Y4B	C07-C03	4.36	1.64	1.56
11	B	502	Y4B	C04-C05	4.22	1.61	1.53
11	B	502	Y4B	C07-C03	3.88	1.63	1.56
10	A	501	PIO	P5-O5	3.18	1.65	1.59
10	C	601	PIO	P4-O4	3.16	1.65	1.59
10	A	501	PIO	P4-O4	3.12	1.65	1.59
10	C	601	PIO	P5-O5	3.06	1.65	1.59
10	C	601	PIO	O3C-C1B	2.46	1.40	1.33
10	A	501	PIO	O3C-C1B	2.43	1.40	1.33
10	C	601	PIO	O2C-C1A	2.35	1.40	1.34
10	A	501	PIO	O2C-C2C	-2.31	1.40	1.46
11	A	502	Y4B	C04-C03	2.26	1.57	1.53
10	A	501	PIO	O2C-C1A	2.26	1.40	1.34
11	B	502	Y4B	C04-C03	2.24	1.57	1.53
10	C	601	PIO	P5-O51	-2.19	1.46	1.54
10	A	501	PIO	P5-O51	-2.19	1.46	1.54
10	C	601	PIO	P4-O41	-2.18	1.46	1.54
10	A	501	PIO	P4-O41	-2.18	1.46	1.54
10	A	501	PIO	P4-O43	-2.17	1.46	1.54
10	C	601	PIO	P4-O43	-2.17	1.46	1.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
10	A	501	PIO	P5-O52	-2.16	1.46	1.54
10	C	601	PIO	P5-O52	-2.14	1.46	1.54
10	A	501	PIO	O3C-C3C	-2.06	1.40	1.45
10	C	601	PIO	O3C-C3C	-2.05	1.40	1.45
10	C	601	PIO	O2C-C2C	-2.02	1.41	1.46
11	A	502	Y4B	C13-C05	2.01	1.58	1.54

All (41) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
11	A	502	Y4B	C23-C22-C12	4.94	124.74	117.56
11	B	502	Y4B	C23-C22-C12	4.84	124.60	117.56
11	A	502	Y4B	C10-C06-C12	4.42	121.49	116.10
11	B	502	Y4B	O02-C22-C23	-4.13	113.82	121.15
11	A	502	Y4B	O02-C22-C23	-4.07	113.92	121.15
11	B	502	Y4B	C10-C06-C12	4.05	121.04	116.10
11	B	502	Y4B	C13-C05-C04	4.03	125.72	119.08
11	A	502	Y4B	C13-C05-C04	3.99	125.64	119.08
10	A	501	PIO	O2C-C1A-C2A	3.97	120.07	111.50
10	C	601	PIO	O2C-C1A-C2A	3.95	120.02	111.50
11	A	502	Y4B	C05-C04-C03	3.81	114.19	109.09
11	A	502	Y4B	C09-C03-C04	-3.80	106.28	111.75
11	B	502	Y4B	C18-C06-C10	-3.56	104.97	110.59
11	B	502	Y4B	C12-C06-C05	3.54	103.50	99.72
11	B	502	Y4B	C09-C03-C04	-3.54	106.66	111.75
11	A	502	Y4B	C13-C05-C06	-3.41	99.73	103.84
11	B	502	Y4B	C07-C03-C04	3.13	115.71	112.42
11	A	502	Y4B	C16-C12-C06	-3.08	101.46	104.21
11	A	502	Y4B	C15-C20-C21	3.01	114.33	110.47
11	A	502	Y4B	C07-C03-C04	2.79	115.35	112.42
11	B	502	Y4B	C14-C11-C04	2.71	116.60	112.14
11	B	502	Y4B	C05-C04-C03	2.69	112.70	109.09
11	B	502	Y4B	C13-C05-C06	-2.69	100.60	103.84
11	A	502	Y4B	C18-C06-C10	-2.59	106.50	110.59
11	B	502	Y4B	C18-C06-C12	-2.56	106.09	110.24
11	A	502	Y4B	C15-C07-C03	2.55	115.36	111.35
10	A	501	PIO	O3C-C1B-C2B	2.50	119.75	111.91
11	B	502	Y4B	C15-C07-C03	2.48	115.25	111.35
10	C	601	PIO	O3C-C1B-C2B	2.41	119.45	111.91
11	B	502	Y4B	C20-C15-C07	2.38	116.86	112.78
11	A	502	Y4B	C14-C11-C04	2.37	116.04	112.14
11	B	502	Y4B	C19-C07-C15	-2.35	104.48	108.26

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
10	C	601	PIO	O11-P1-O12	-2.31	100.83	112.24
11	A	502	Y4B	C20-C15-C07	2.24	116.62	112.78
11	B	502	Y4B	C14-C08-C07	-2.19	108.37	112.31
11	A	502	Y4B	C08-C17-C21	2.19	115.97	112.76
11	B	502	Y4B	C15-C20-C21	2.18	113.26	110.47
10	A	501	PIO	O11-P1-O12	-2.17	101.49	112.24
11	A	502	Y4B	C11-C04-C03	-2.10	107.89	110.49
10	A	501	PIO	O1-C1-C2	2.09	113.52	108.66
12	E	501	ABU	CB-CG-C	-2.06	109.28	114.47

There are no chirality outliers.

All (33) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
10	A	501	PIO	C2-C1-O1-P1
10	C	601	PIO	C2A-C1A-O2C-C2C
10	C	601	PIO	O1B-C1B-O3C-C3C
10	C	601	PIO	O1A-C1A-O2C-C2C
10	C	601	PIO	C2B-C1B-O3C-C3C
10	C	601	PIO	C1B-C2B-C3B-C4B
10	A	501	PIO	C1B-C2B-C3B-C4B
10	A	501	PIO	C2B-C1B-O3C-C3C
10	C	601	PIO	C1C-O13-P1-O1
10	C	601	PIO	C4A-C5A-C6A-C7A
10	A	501	PIO	O1B-C1B-O3C-C3C
10	A	501	PIO	C1A-C2A-C3A-C4A
10	A	501	PIO	C4A-C5A-C6A-C7A
10	C	601	PIO	C4B-C5B-C6B-C7B
10	A	501	PIO	C3B-C4B-C5B-C6B
10	C	601	PIO	C1C-C2C-C3C-O3C
10	C	601	PIO	C1-O1-P1-O13
10	A	501	PIO	C4B-C5B-C6B-C7B
10	A	501	PIO	C4-O4-P4-O42
10	C	601	PIO	C1C-C2C-O2C-C1A
10	A	501	PIO	O2C-C2C-C3C-O3C
10	C	601	PIO	O2C-C2C-C3C-O3C
10	C	601	PIO	C1C-O13-P1-O11
10	A	501	PIO	C5A-C6A-C7A-C8A
10	A	501	PIO	C1C-O13-P1-O1
10	C	601	PIO	C5B-C6B-C7B-C8B
10	C	601	PIO	C5A-C6A-C7A-C8A
12	E	501	ABU	OXT-C-CG-CB

Continued on next page...

Continued from previous page...

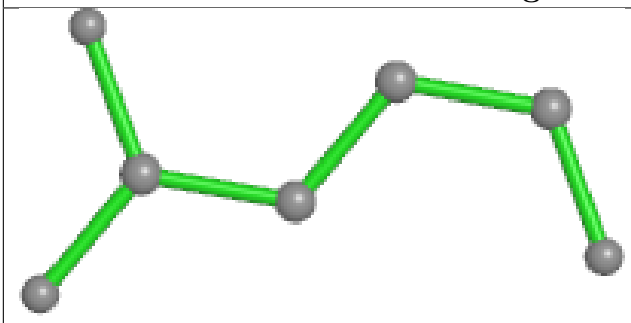
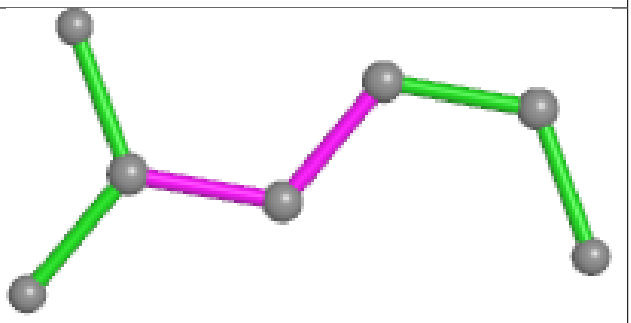
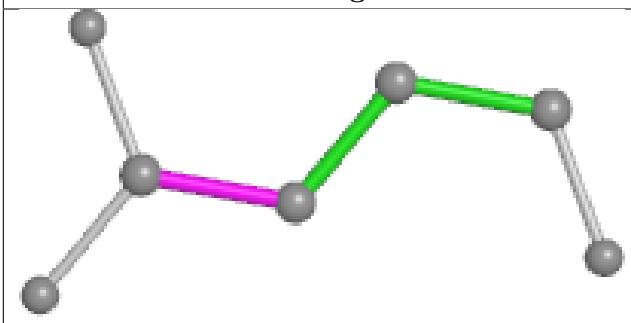
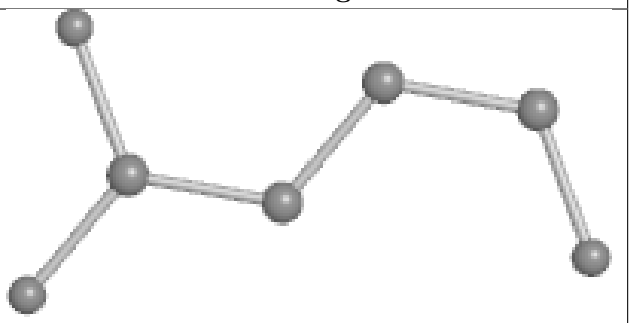
Mol	Chain	Res	Type	Atoms
12	E	501	ABU	O-C-CG-CB
12	B	501	ABU	O-C-CG-CB
12	B	501	ABU	OXT-C-CG-CB
10	C	601	PIO	C4-O4-P4-O42
10	A	501	PIO	C1C-C2C-C3C-O3C

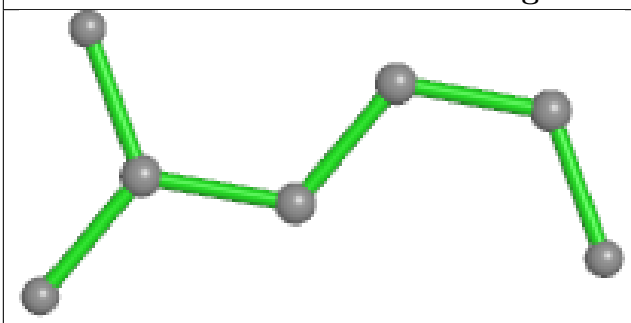
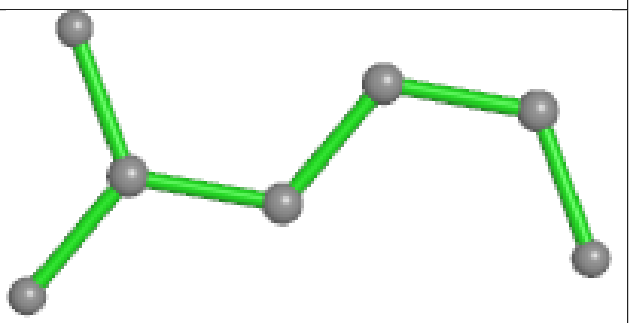
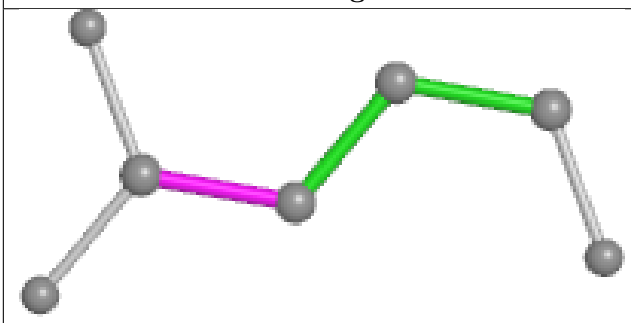
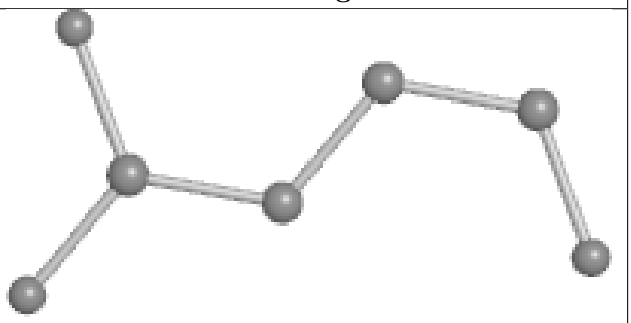
There are no ring outliers.

3 monomers are involved in 5 short contacts:

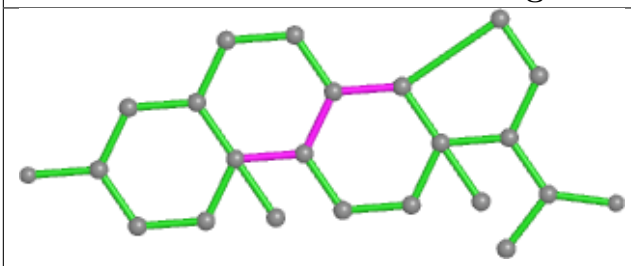
Mol	Chain	Res	Type	Clashes	Symm-Clashes
12	B	501	ABU	1	0
11	B	502	Y4B	1	0
10	A	501	PIO	3	0

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.

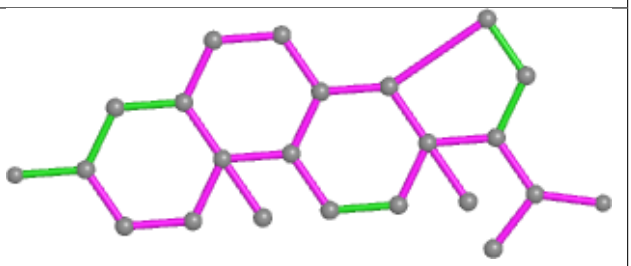
Ligand ABU E 501	
 Bond lengths	 Bond angles
 Torsions	 Rings

Ligand ABU B 501	
 Bond lengths	 Bond angles
 Torsions	 Rings

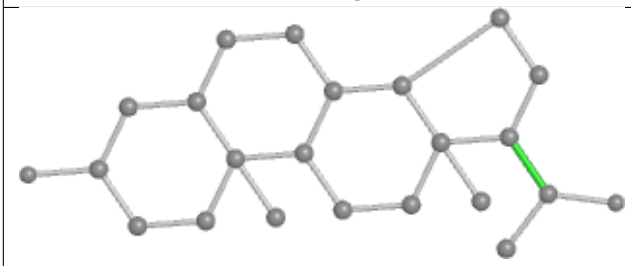
Ligand Y4B B 502



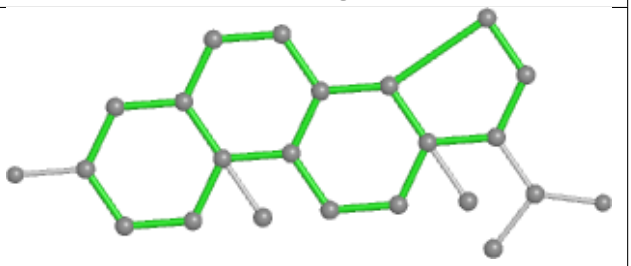
Bond lengths



Bond angles

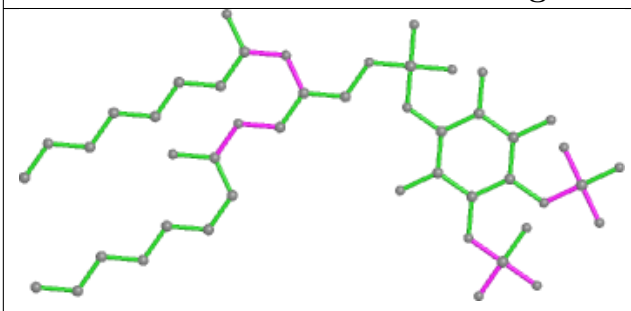


Torsions

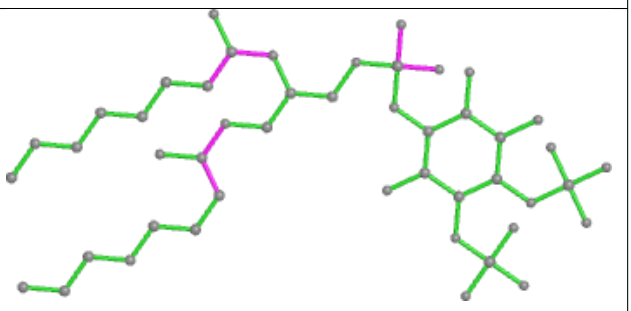


Rings

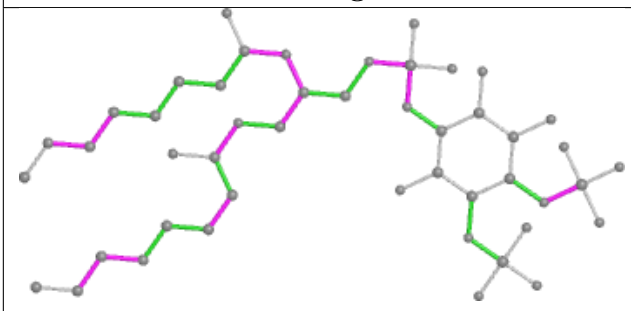
Ligand PIO C 601



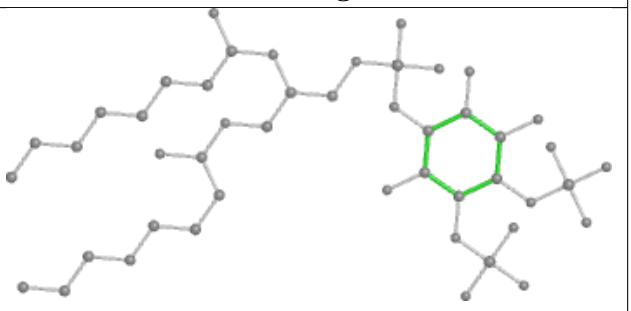
Bond lengths



Bond angles

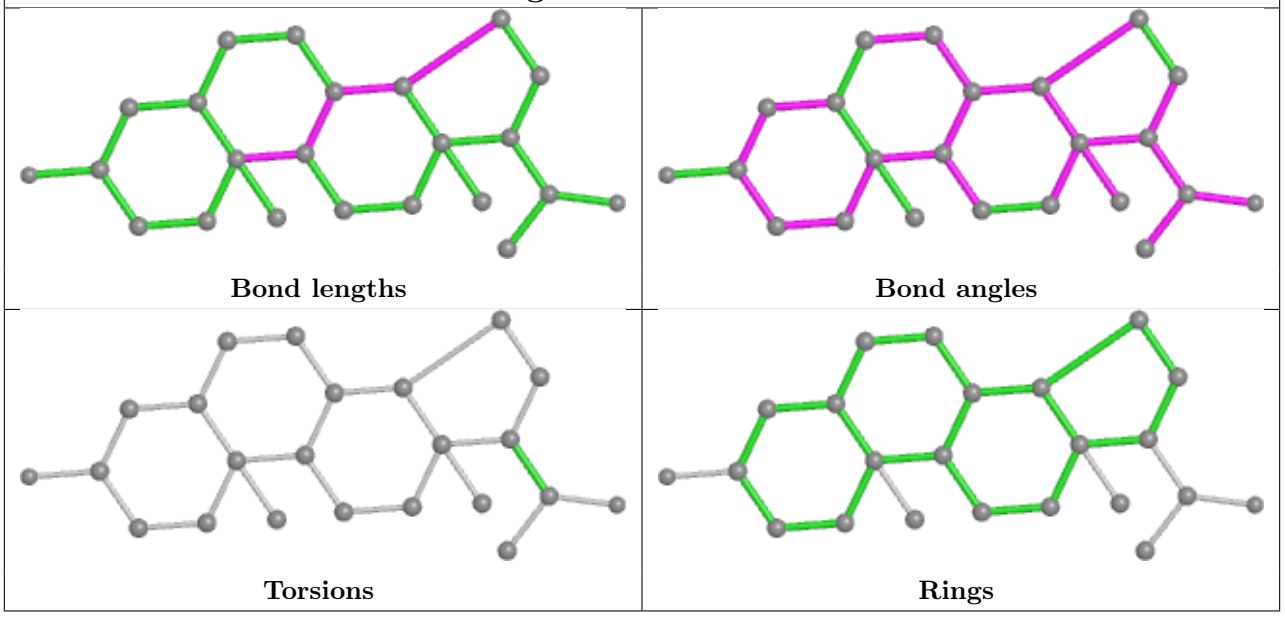


Torsions

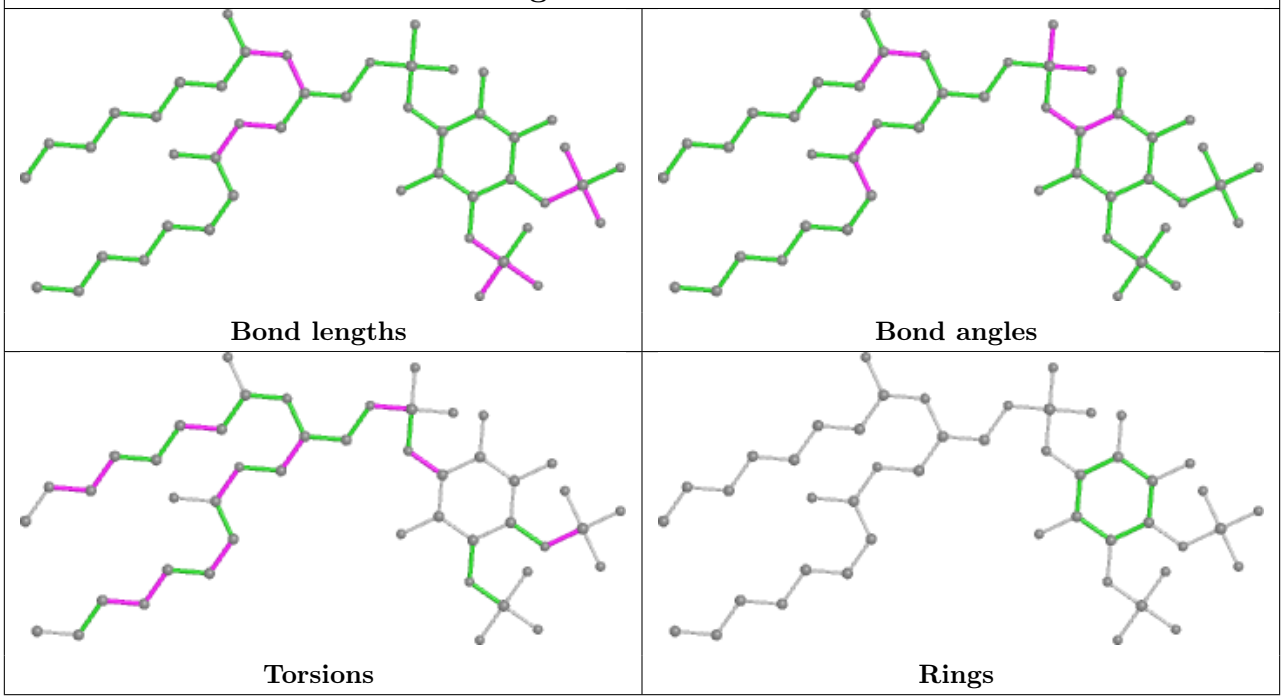


Rings

Ligand Y4B A 502



Ligand PIO A 501



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.

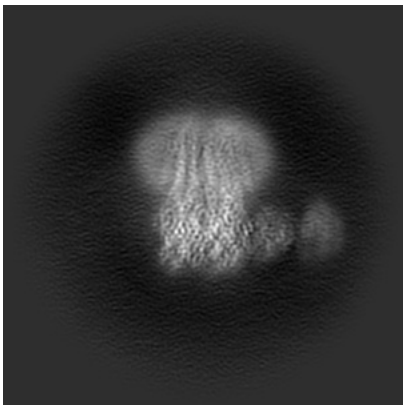
6 Map visualisation

This section contains visualisations of the EMDB entry EMD-29741. These allow visual inspection of the internal detail of the map and identification of artifacts.

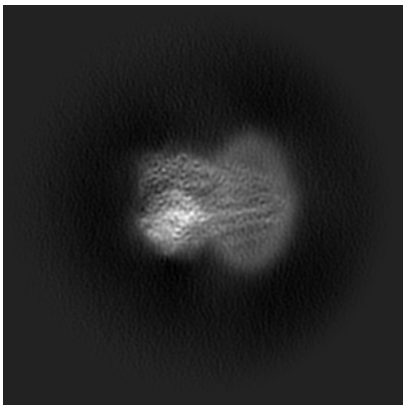
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections

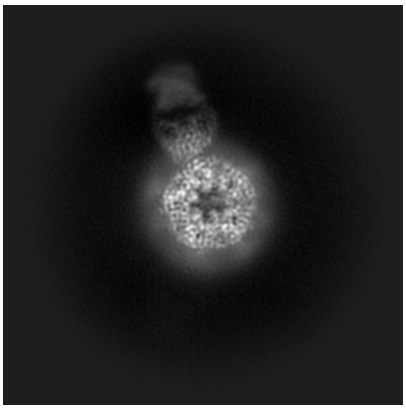
6.1.1 Primary map



X

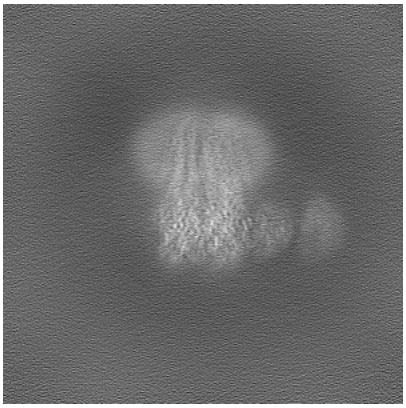


Y

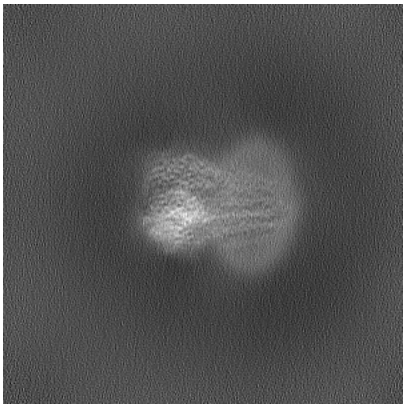


Z

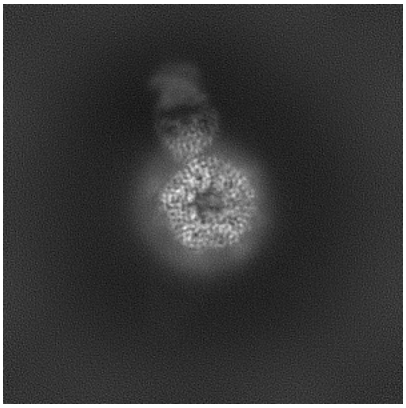
6.1.2 Raw map



X



Y

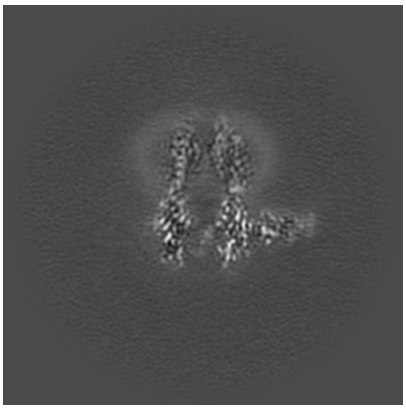


Z

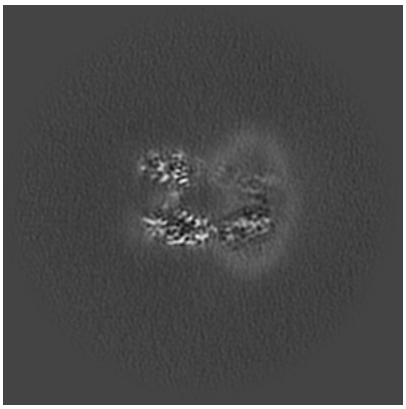
The images above show the map projected in three orthogonal directions.

6.2 Central slices (i)

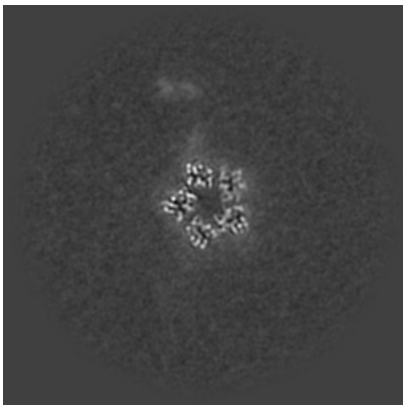
6.2.1 Primary map



X Index: 180

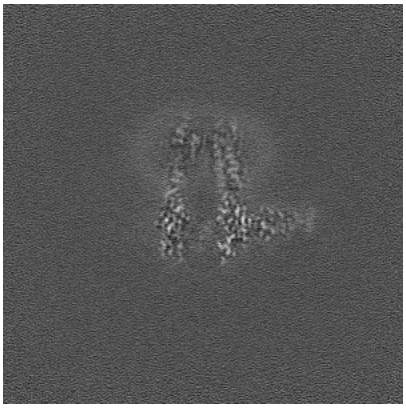


Y Index: 180

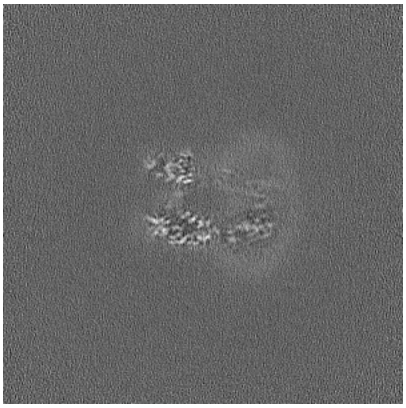


Z Index: 180

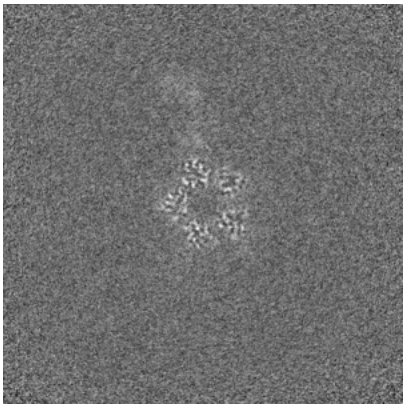
6.2.2 Raw map



X Index: 180



Y Index: 180

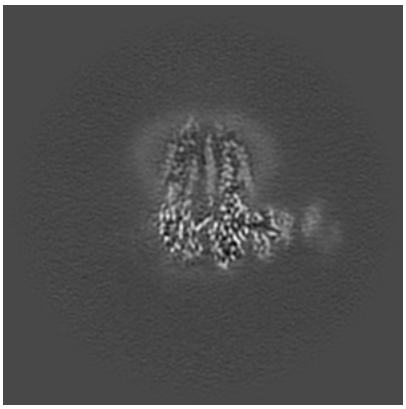


Z Index: 180

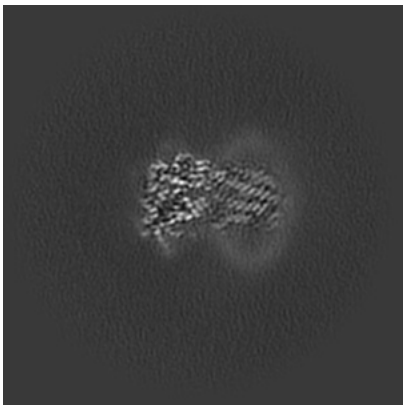
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices (i)

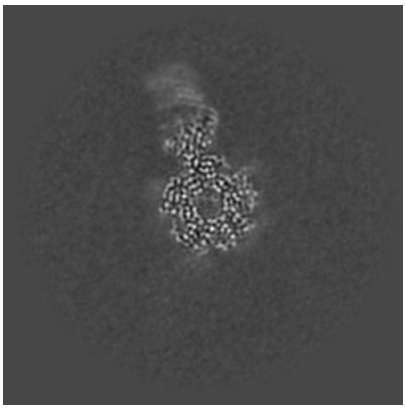
6.3.1 Primary map



X Index: 172

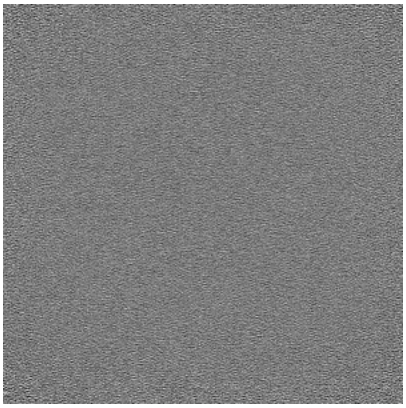


Y Index: 161

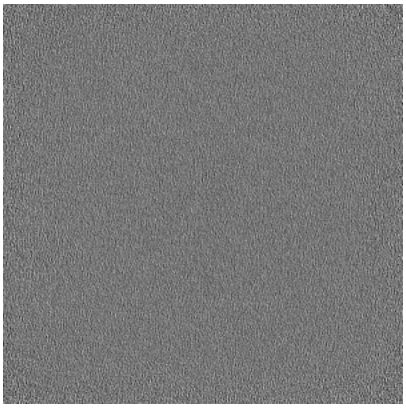


Z Index: 159

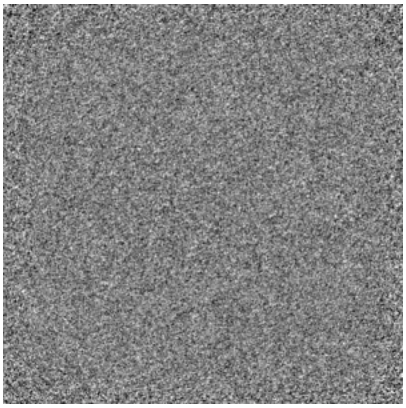
6.3.2 Raw map



X Index: 0



Y Index: 0

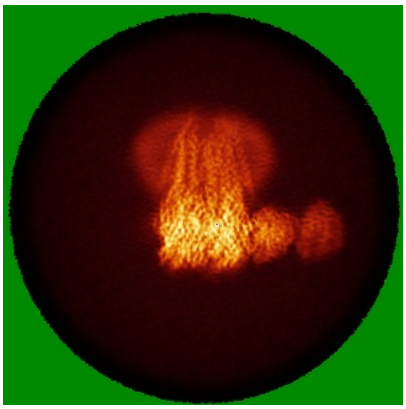


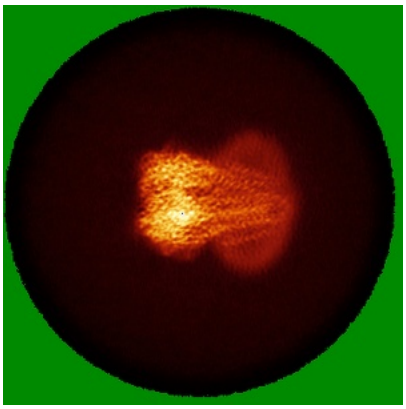
Z Index: 0

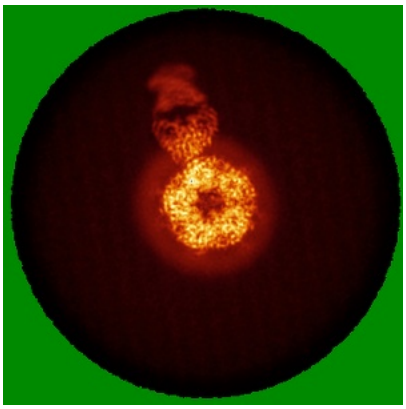
The images above show the largest variance slices of the map in three orthogonal directions.

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

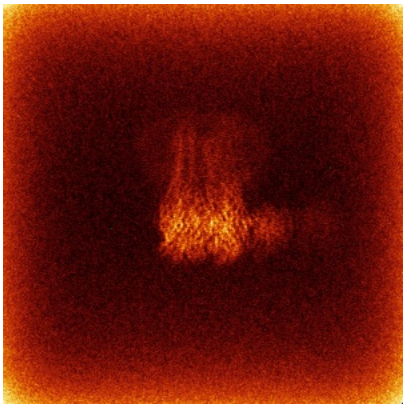
6.4.1 Primary map

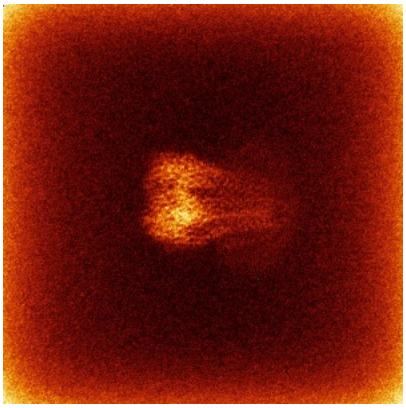

X

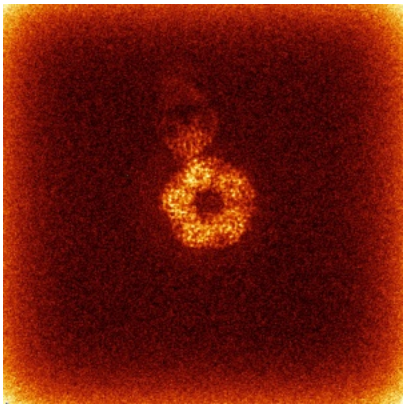

Y


Z

6.4.2 Raw map


X

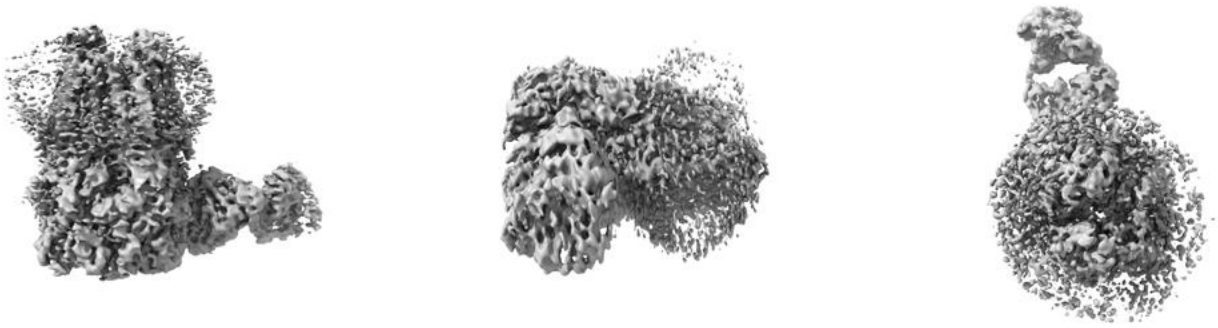

Y


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 0.09. 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.

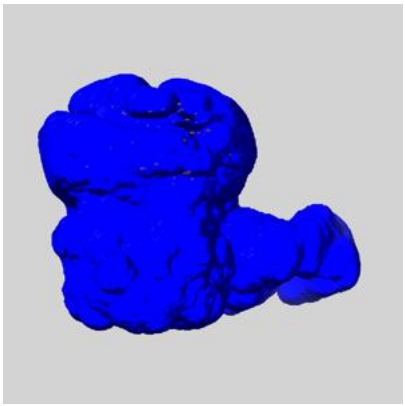
6.6 Mask visualisation

This section shows the 3D surface view of the primary map at 50% transparency overlaid with the specified mask at 0% transparency

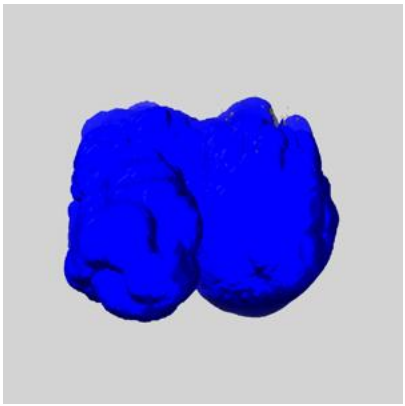
A mask typically either:

- Encompasses the whole structure
- Separates out a domain, a functional unit, a monomer or an area of interest from a larger structure

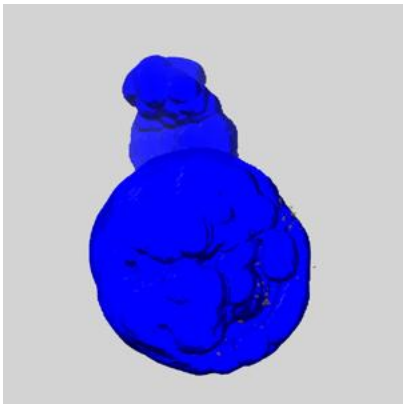
6.6.1 emd_29741_msk_1.map



X



Y

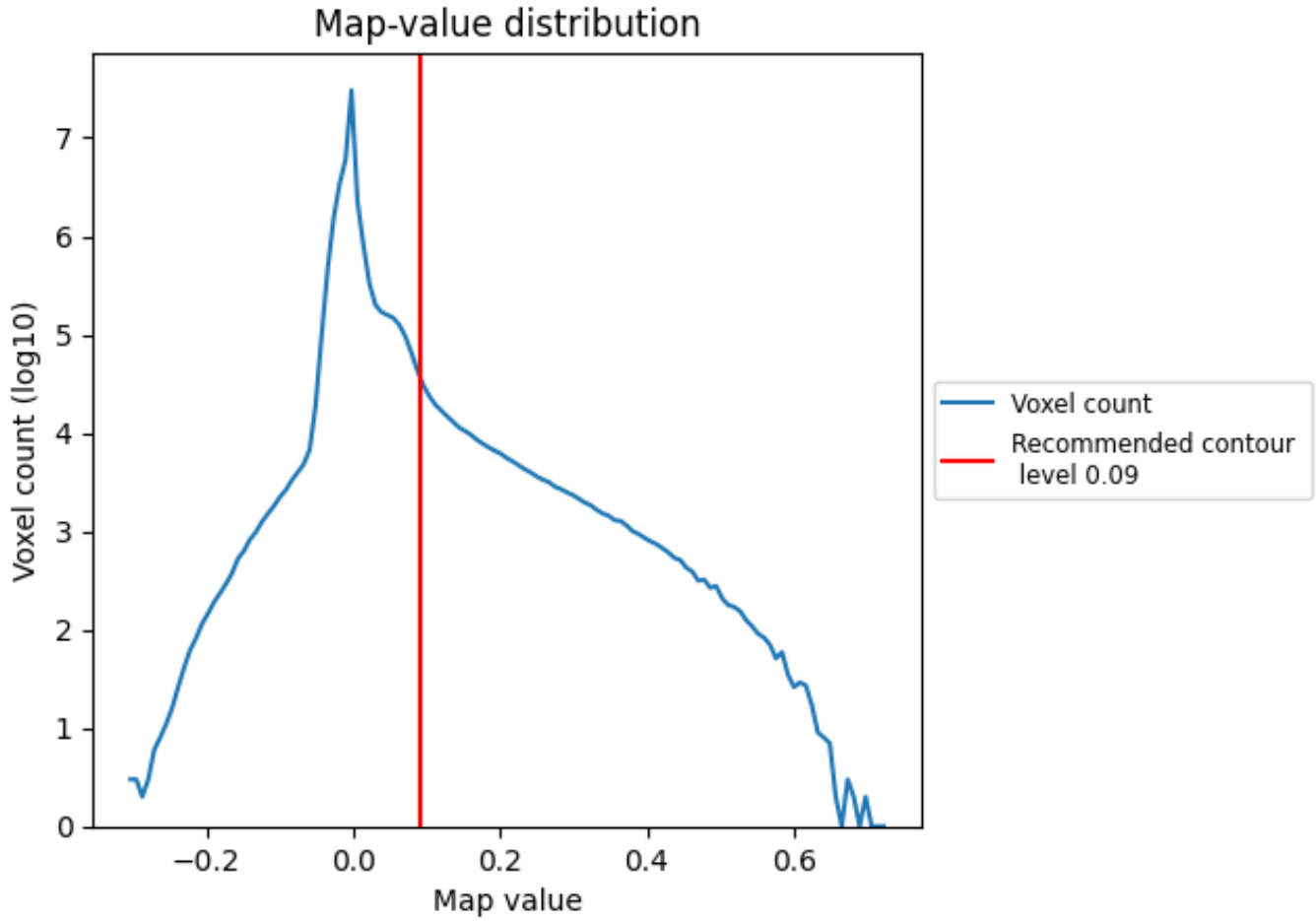


Z

7 Map analysis

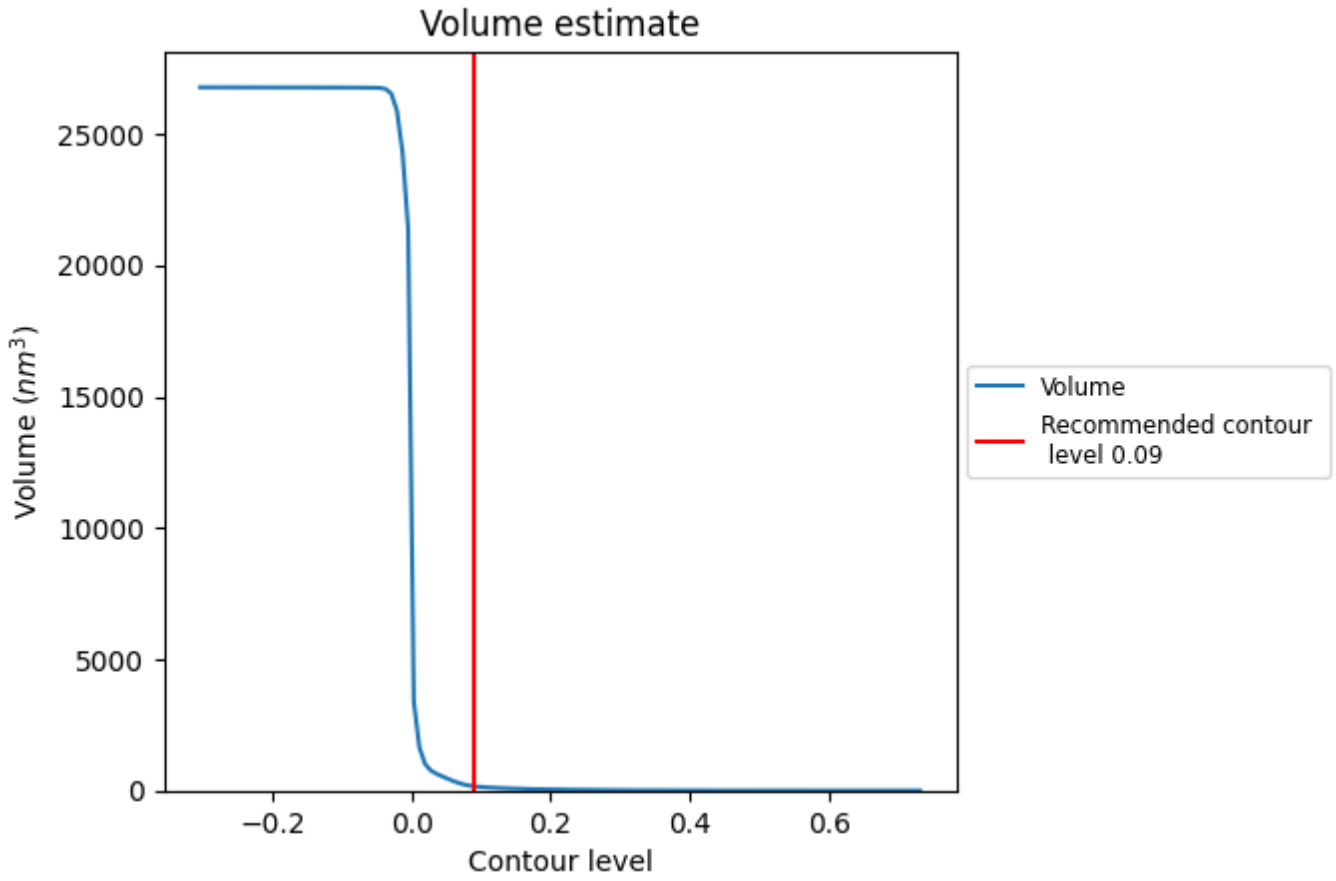
This section contains the results of statistical analysis of the map.

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.

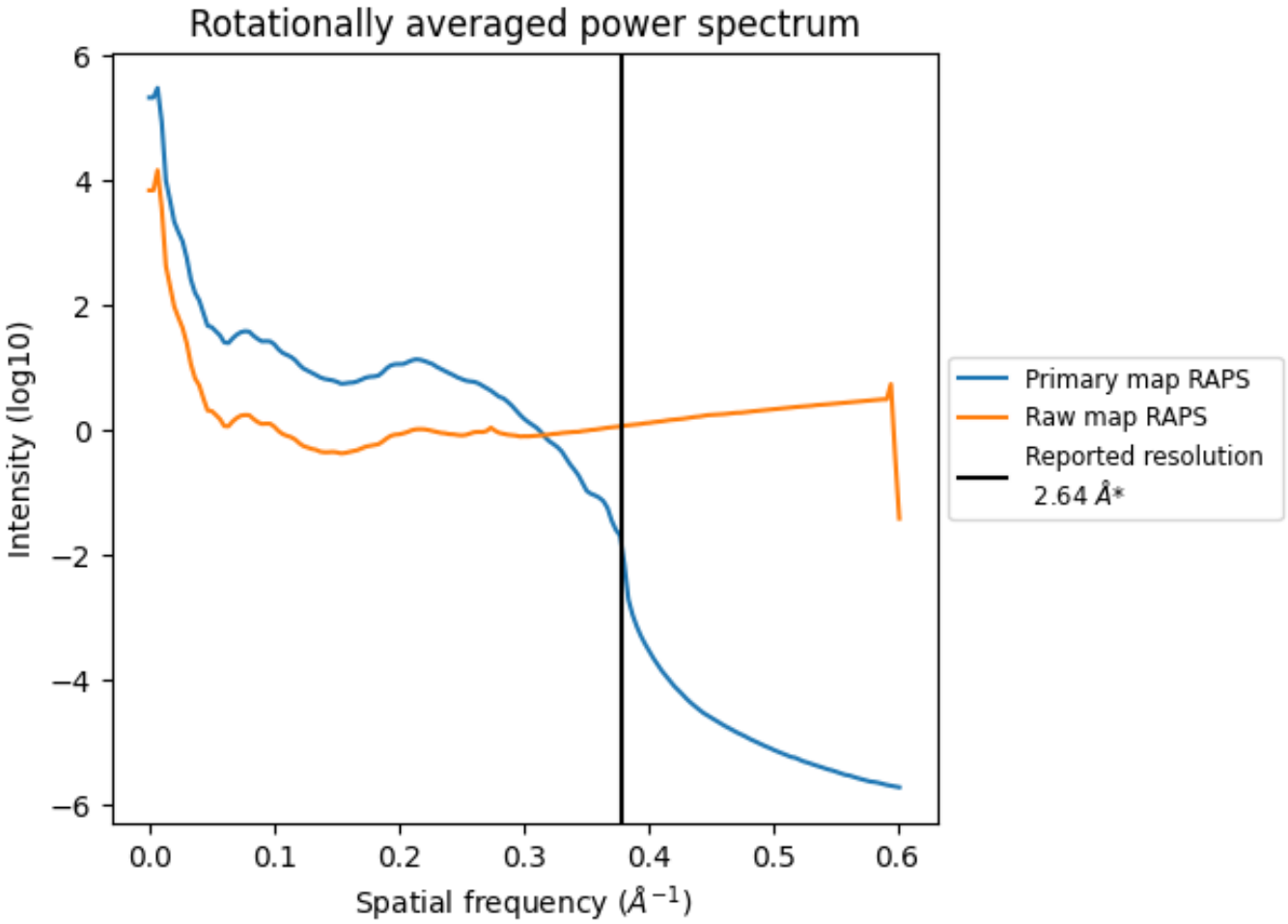
7.2 Volume estimate



The volume at the recommended contour level is 163 nm^3; this corresponds to an approximate mass of 147 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)

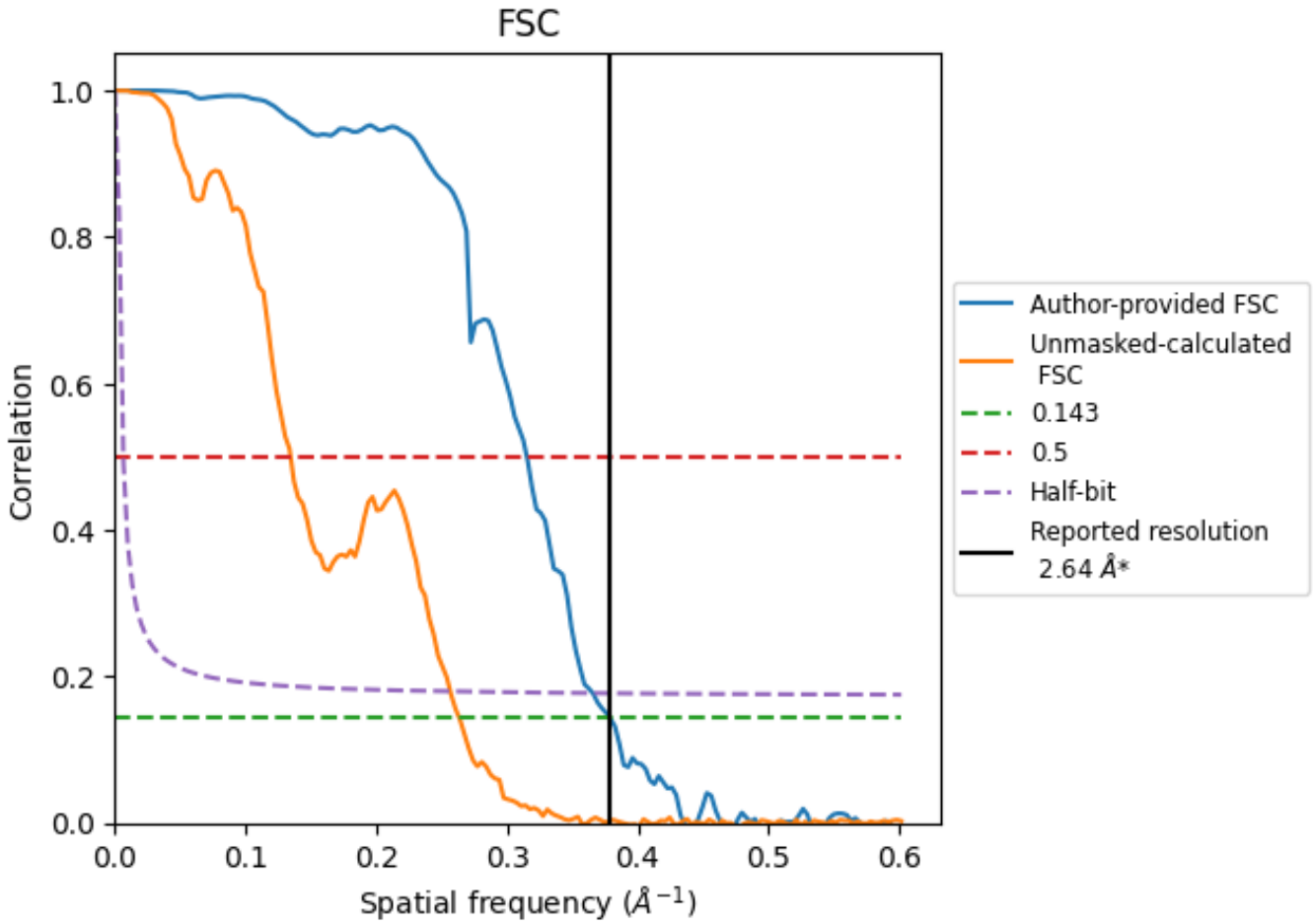


*Reported resolution corresponds to spatial frequency of 0.379 Å^{-1}

8 Fourier-Shell correlation

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



*Reported resolution corresponds to spatial frequency of 0.379 Å^{-1}

8.2 Resolution estimates (i)

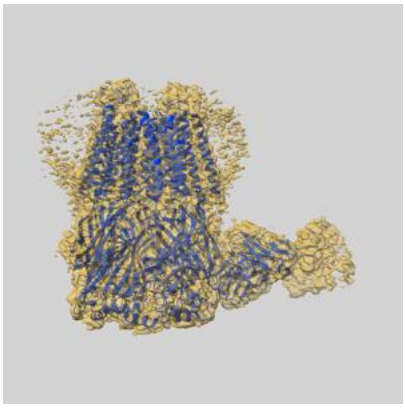
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.64	-	-
Author-provided FSC curve	2.64	3.17	2.73
Unmasked-calculated*	3.79	7.43	3.90

*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 3.79 differs from the reported value 2.64 by more than 10 %

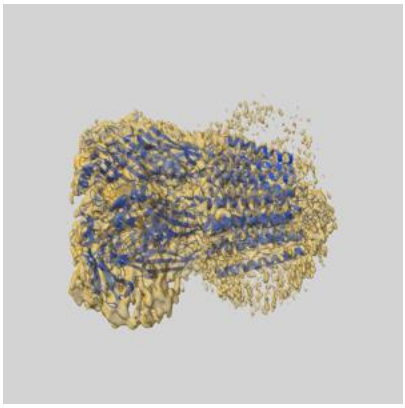
9 Map-model fit

This section contains information regarding the fit between EMDB map EMD-29741 and PDB model 8G5F. Per-residue inclusion information can be found in section 3 on page 9.

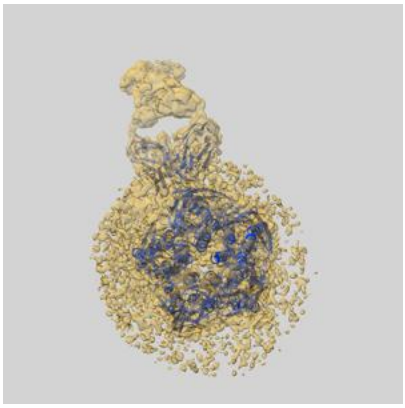
9.1 Map-model overlay



X



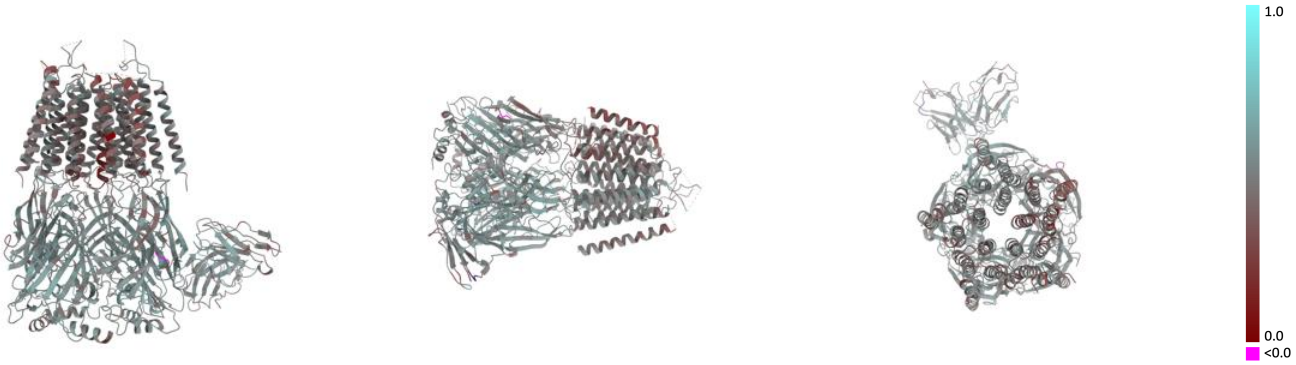
Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.09 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.

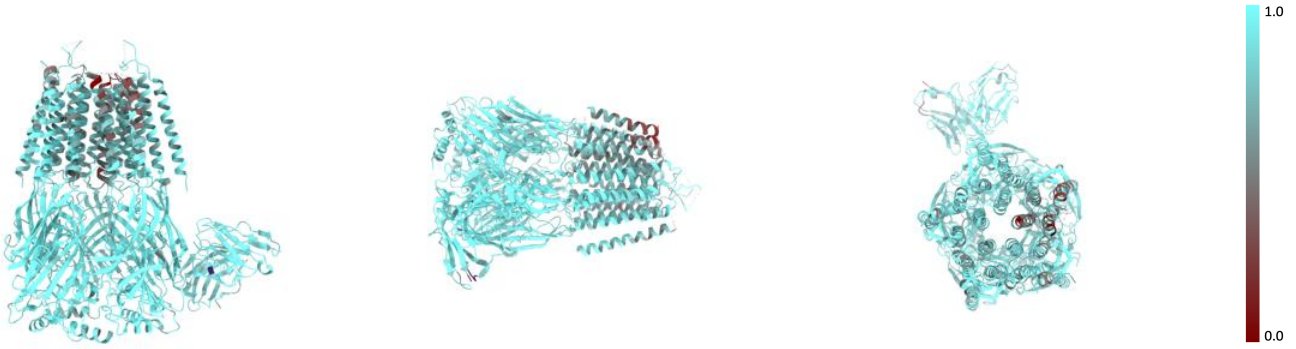
9.2 Q-score mapped to coordinate model (i)



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.

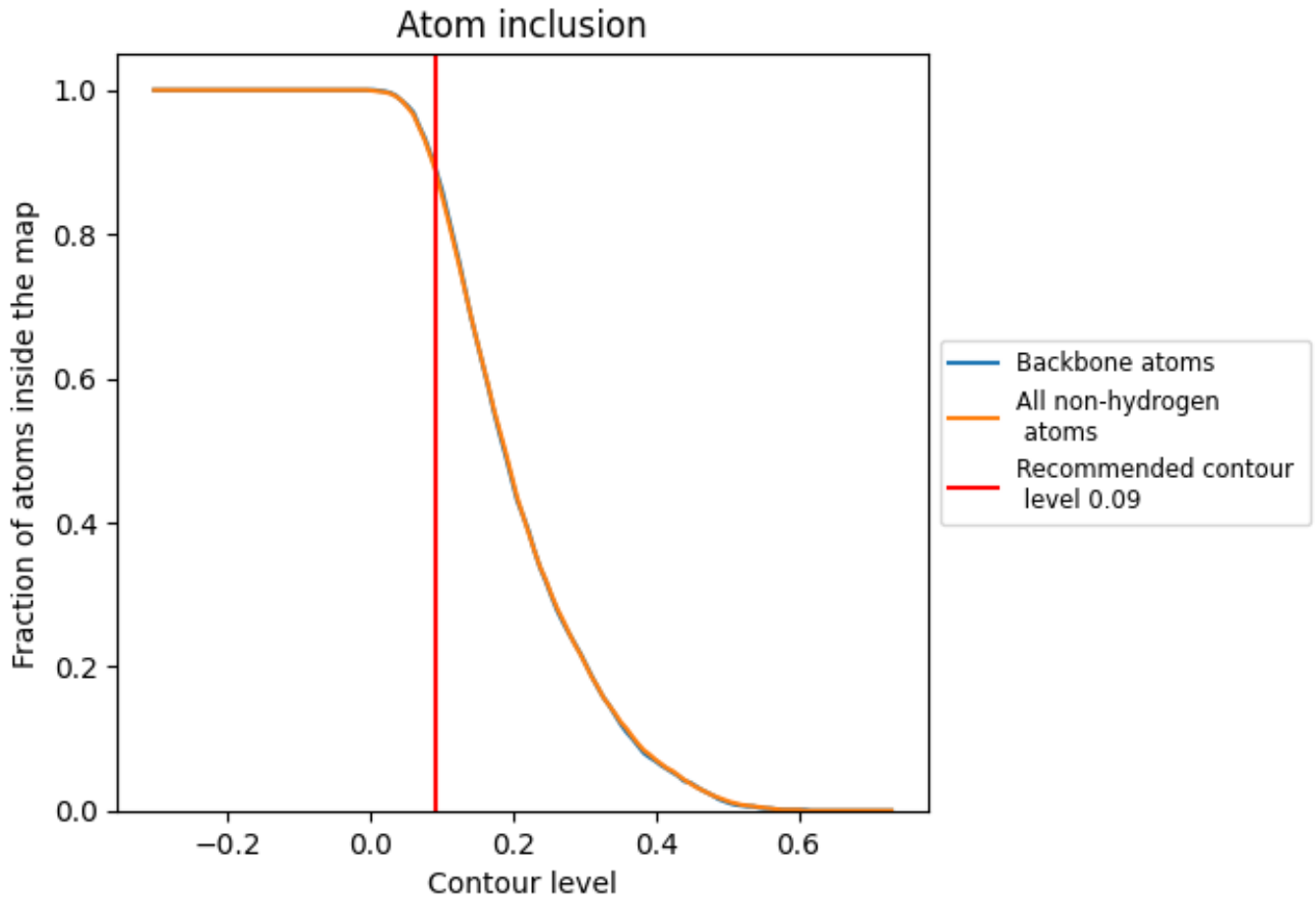
9.3 Atom inclusion mapped to coordinate model (i)



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 (0.09).

9.4 Atom inclusion (i)



At the recommended contour level, 90% of all backbone atoms, 89% 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 (0.09) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.8920	0.4910
A	0.8900	0.4850
B	0.9160	0.5030
C	0.9270	0.5240
D	0.8280	0.4520
E	0.9040	0.4960
F	0.8290	0.4070
G	0.7440	0.4310
H	0.8970	0.4040
I	0.8210	0.4580
J	0.8980	0.4900
K	0.9290	0.5020
L	0.3210	0.2370
M	0.5640	0.2780
N	0.8970	0.3980