Full wwPDB EM Validation Report (i)

Nov 20, 2022 – 11:27 am GMT

PDB ID	:	6QNO
EMDB ID	:	EMD-4598
Title	:	Rhodopsin-Gi protein complex
Authors	:	Tsai, C.-J.; Marino, J.; Adaixo, R.J.; Pamula, F.; Muehle, J.; Maeda, S.; Flock, T.; Taylor, N.M.I.; Mohammed, I.; Matile, H.; Dawson, R.J.P.; Deupi, X.; Stahlberg, H.; Schertler, G.F.X.
Deposited on	:	2019-02-11
Resolution	:	4.38 Å(reported)
Based on initial models	:	6FUF, 1GOT, 6QNK

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

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

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

The following versions of software and data (see references (1)) were used in the production of this report:

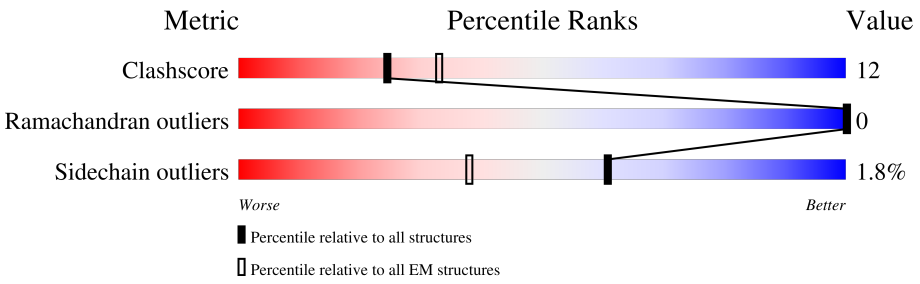
EMDB validation analysis	:	0.0.1.dev43
Mogul	:	1.8.4, CSD as541be (2020)
MolProbity	:	4.02b-467
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
MapQ	:	1.9.9
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.31.2

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 4.38 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	376	30%; 38%; 17%; •; 44%
2	B	340	•; 58%; 32%; •; 8%
3	G	74	14%; 38%; 11%; 51%
4	L	239	5%; 62%; 28%; •; 9%
5	H	249	8%; 63%; 25%; 11%
6	R	348	17%; 77%; 18%; •; •
7	C	2	100%; 100%

2 Entry composition (i)

There are 8 unique types of molecules in this entry. The entry contains 10429 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 Guanine nucleotide-binding protein G(i) subunit alpha-1.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	A	212	1711	1089	286	324	12	0	0

There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-21	MET	-	initiating methionine	UNP P63096
A	-20	LYS	-	expression tag	UNP P63096
A	-19	LYS	-	expression tag	UNP P63096
A	-18	HIS	-	expression tag	UNP P63096
A	-17	HIS	-	expression tag	UNP P63096
A	-16	HIS	-	expression tag	UNP P63096
A	-15	HIS	-	expression tag	UNP P63096
A	-14	HIS	-	expression tag	UNP P63096
A	-13	HIS	-	expression tag	UNP P63096
A	-12	HIS	-	expression tag	UNP P63096
A	-11	HIS	-	expression tag	UNP P63096
A	-10	HIS	-	expression tag	UNP P63096
A	-9	HIS	-	expression tag	UNP P63096
A	-8	GLU	-	expression tag	UNP P63096
A	-7	ASN	-	expression tag	UNP P63096
A	-6	LEU	-	expression tag	UNP P63096
A	-5	TYR	-	expression tag	UNP P63096
A	-4	PHE	-	expression tag	UNP P63096
A	-3	GLN	-	expression tag	UNP P63096
A	-2	GLY	-	expression tag	UNP P63096
A	-1	GLY	-	expression tag	UNP P63096
A	0	SER	-	expression tag	UNP P63096

- Molecule 2 is a protein called Guanine nucleotide-binding protein G(I)/G(S)/G(T) subunit beta-1.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
2	B	312	2392	1481	425	466	20	0	0

- Molecule 3 is a protein called Guanine nucleotide-binding protein G(T) subunit gamma-T1.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
3	G	36	293	183	47	60	3	0	0

- Molecule 4 is a protein called Fab antibody fragment light chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
4	L	217	1679	1049	284	338	8	0	0

- Molecule 5 is a protein called Fab antibody fragment heavy chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
5	H	221	1660	1051	267	333	9	2	0

- Molecule 6 is a protein called Rhodopsin.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
6	R	334	2646	1758	407	455	26	0	0

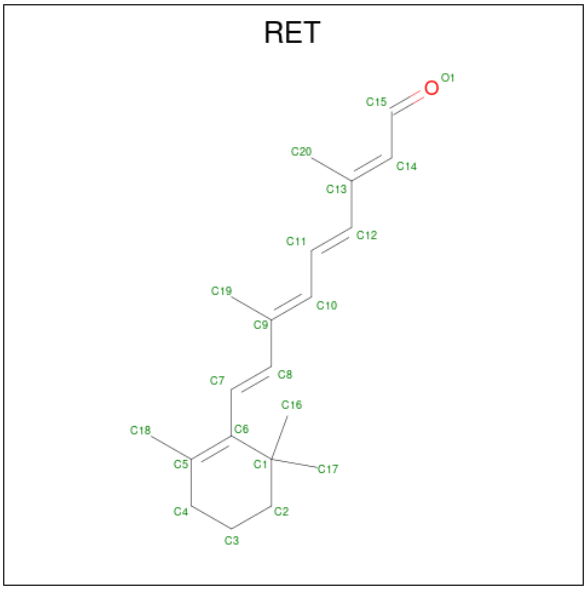
There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
R	2	CYS	ASN	engineered mutation	UNP P02699
R	257	TYR	MET	engineered mutation	UNP P02699
R	282	CYS	ASP	engineered mutation	UNP P02699

- Molecule 7 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
			Total	C	N	O		
7	C	2	28	16	2	10	0	0

- Molecule 8 is RETINAL (three-letter code: RET) (formula: $C_{20}H_{28}O$) (labeled as "Ligand of Interest" by depositor).



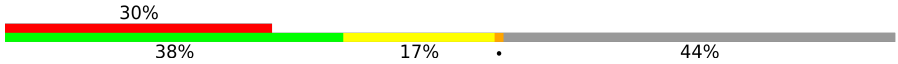
Mol	Chain	Residues	Atoms		AltConf
			Total	C	
8	R	1	20	20	0

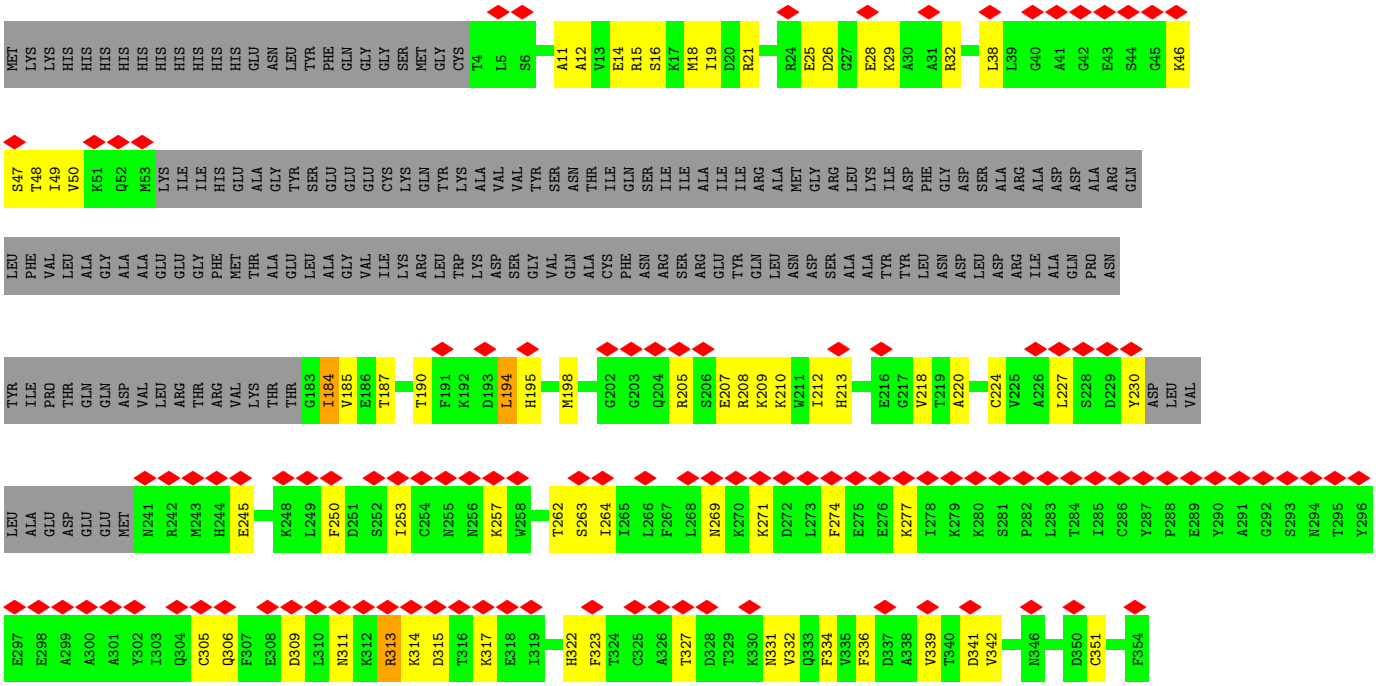
3 Residue-property plots (i)

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

• Molecule 1: Guanine nucleotide-binding protein G(i) subunit alpha-1

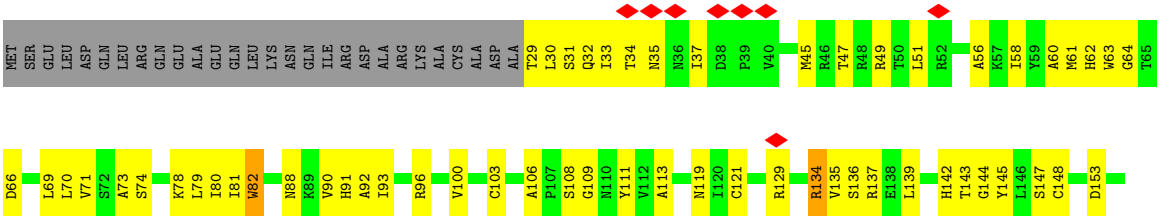
Chain A:

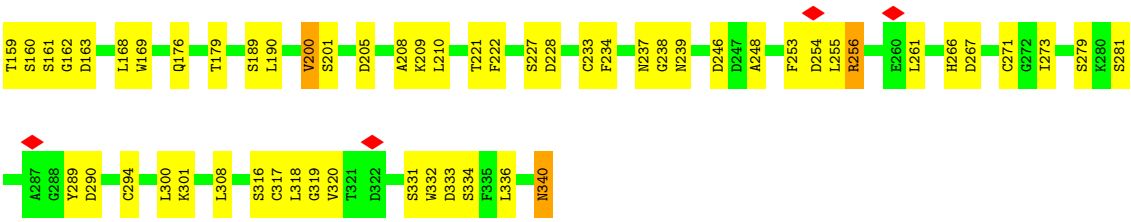




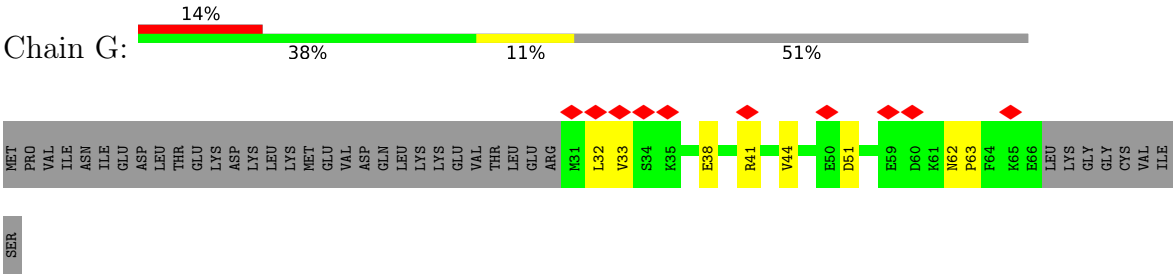
• Molecule 2: Guanine nucleotide-binding protein G(I)/G(S)/G(T) subunit beta-1

Chain B:

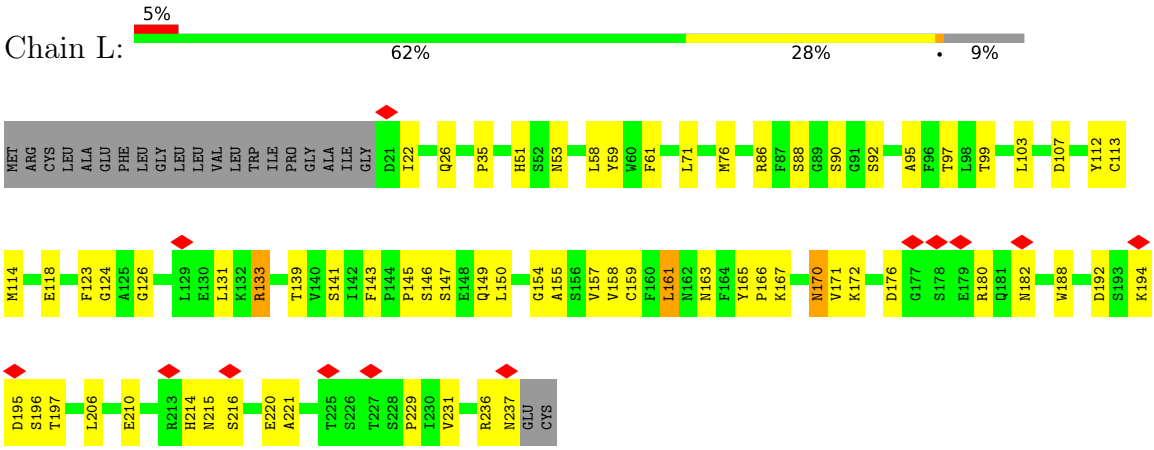




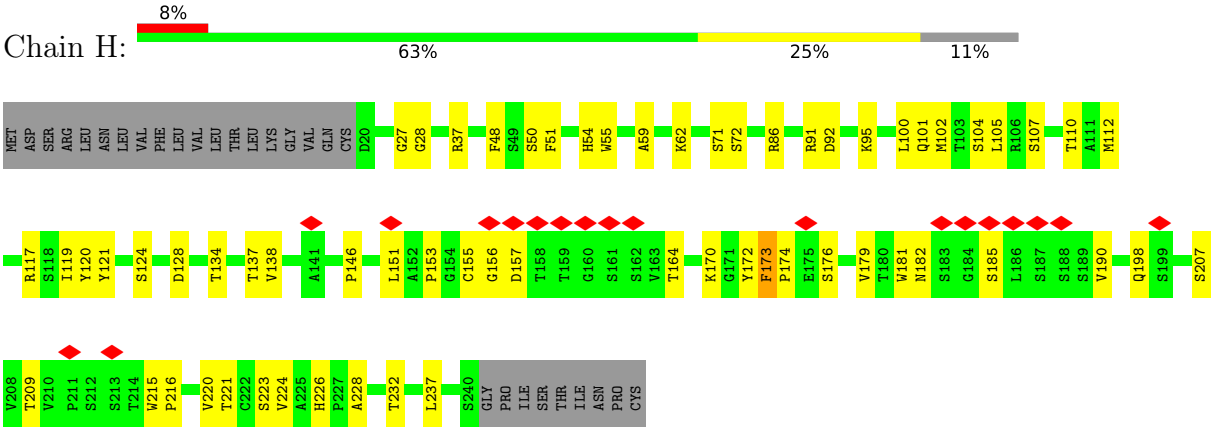
• Molecule 3: Guanine nucleotide-binding protein G(T) subunit gamma-T1



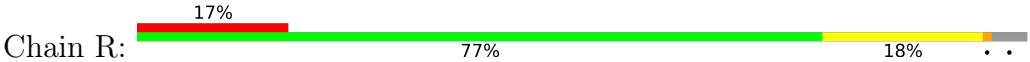
• Molecule 4: Fab antibody fragment light chain

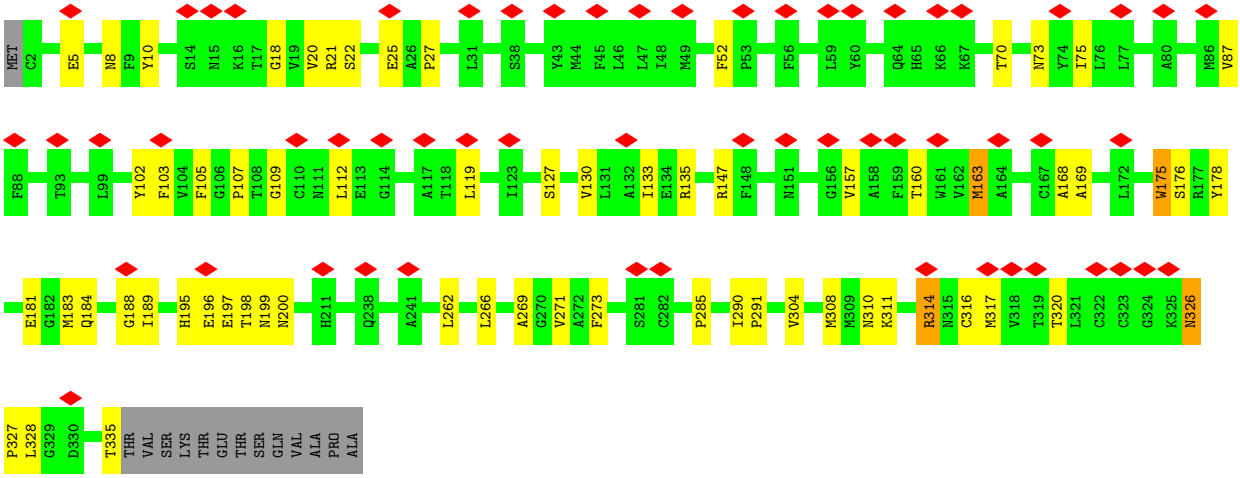


• Molecule 5: Fab antibody fragment heavy chain



• Molecule 6: Rhodopsin





• Molecule 7: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain C:

4 Experimental information (i)

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	115000	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 ($e^-/Å^2$)	60	Depositor
Minimum defocus (nm)	15000	Depositor
Maximum defocus (nm)	25000	Depositor
Magnification	165000	Depositor
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	0.120	Depositor
Minimum map value	-0.051	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.003	Depositor
Recommended contour level	0.03	Depositor
Map size (Å)	280.54, 280.54, 280.54	wwPDB
Map dimensions	338, 338, 338	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.83000004, 0.83000004, 0.83000004	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: NAG, RET

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	0.39	0/1740	0.69	1/2331 (0.0%)
2	B	0.59	0/2439	0.74	0/3310
3	G	0.44	0/298	0.71	0/398
4	L	0.50	0/1717	0.67	1/2333 (0.0%)
5	H	0.53	0/1712	0.68	0/2337
6	R	0.46	0/2730	0.64	0/3722
All	All	0.50	0/10636	0.68	2/14431 (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
5	H	0	1
6	R	0	3
All	All	0	4

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	341	ASP	CB-CG-OD1	8.20	125.68	118.30
4	L	123	PHE	CB-CG-CD1	6.55	125.38	120.80

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
5	H	173	PHE	Peptide
6	R	169	ALA	Peptide
6	R	175	TRP	Peptide
6	R	327	PRO	Peptide

5.2 Too-close contacts (i)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1711	0	1700	43	0
2	B	2392	0	2298	80	0
3	G	293	0	280	9	0
4	L	1679	0	1620	41	0
5	H	1660	0	1600	40	0
6	R	2646	0	2605	45	0
7	C	28	0	25	1	0
8	R	20	0	27	1	0
All	All	10429	0	10155	243	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 12.

All (243) 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)
2:B:208:ALA:O	2:B:222:PHE:HB2	1.88	0.73
2:B:45:MET:HA	2:B:340:ASN:O	1.90	0.70
5:H:224:VAL:O	5:H:232:THR:HA	1.94	0.68
2:B:51:LEU:HD22	2:B:336:LEU:HD12	1.76	0.68
1:A:227:LEU:HD13	1:A:274:PHE:HB2	1.76	0.67
5:H:182:ASN:HD21	5:H:220:VAL:HA	1.58	0.67
5:H:190:VAL:HA	5:H:207:SER:O	1.94	0.67
5:H:27:GLY:O	5:H:37:ARG:NH1	2.28	0.66
2:B:33:ILE:O	3:G:41:ARG:NH2	2.28	0.66
6:R:310:ASN:O	6:R:314:ARG:N	2.29	0.66
1:A:224:CYS:SG	1:A:269:ASN:ND2	2.69	0.65
1:A:15:ARG:NH2	5:H:50:SER:O	2.29	0.65

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:L:133:ARG:HH21	4:L:165:TYR:HB2	1.62	0.65
2:B:271:CYS:HB2	2:B:290:ASP:HB2	1.79	0.65
5:H:48:PHE:O	5:H:91:ARG:NH2	2.30	0.64
5:H:101:GLN:NE2	5:H:102:MET:O	2.30	0.64
2:B:121:CYS:HB3	2:B:139:LEU:HB2	1.80	0.64
4:L:172:LYS:HB2	4:L:220:GLU:HB2	1.80	0.63
6:R:27:PRO:HA	6:R:102:TYR:HB3	1.80	0.63
6:R:269:ALA:O	6:R:273:PHE:HB2	2.00	0.62
2:B:51:LEU:HB2	2:B:336:LEU:HB2	1.83	0.61
4:L:59:TYR:HB2	4:L:114:MET:O	2.01	0.61
2:B:69:LEU:HA	2:B:82:TRP:O	2.01	0.60
4:L:145:PRO:HG2	4:L:155:ALA:HB1	1.82	0.60
2:B:233:CYS:SG	2:B:234:PHE:N	2.74	0.60
2:B:64:GLY:N	2:B:69:LEU:O	2.35	0.59
2:B:279:SER:HA	2:B:320:VAL:HG11	1.83	0.59
5:H:155:CYS:SG	5:H:156:GLY:N	2.74	0.59
2:B:159:THR:OG1	2:B:169:TRP:NE1	2.33	0.59
5:H:172:TYR:HD2	5:H:174:PRO:HD2	1.68	0.59
6:R:262:LEU:HD12	6:R:266:LEU:HD22	1.85	0.59
1:A:257:LYS:HZ2	6:R:335:THR:H	1.50	0.58
2:B:32:GLN:O	2:B:35:ASN:ND2	2.36	0.58
2:B:79:LEU:HB3	2:B:93:ILE:HB	1.85	0.58
2:B:161:SER:OG	2:B:162:GLY:N	2.36	0.58
2:B:253:PHE:HA	2:B:261:LEU:HD13	1.86	0.58
2:B:281:SER:HB3	3:G:51:ASP:HB2	1.85	0.57
4:L:176:ASP:HA	4:L:216:SER:HB3	1.86	0.57
2:B:201:SER:O	2:B:208:ALA:HA	2.04	0.57
6:R:10:TYR:O	6:R:184:GLN:NE2	2.38	0.57
2:B:331:SER:OG	2:B:332:TRP:N	2.38	0.57
1:A:187:THR:HG22	1:A:198:MET:HB3	1.87	0.56
5:H:71:SER:OG	5:H:72:SER:N	2.38	0.56
2:B:73:ALA:HB1	2:B:100:VAL:HG11	1.87	0.56
1:A:250:PHE:HA	1:A:253:ILE:HG22	1.88	0.56
4:L:145:PRO:HD3	4:L:157:VAL:HG22	1.86	0.56
1:A:184:ILE:HD12	1:A:185:VAL:H	1.71	0.56
2:B:60:ALA:HB3	2:B:73:ALA:HB3	1.88	0.56
2:B:176:GLN:NE2	2:B:179:THR:OG1	2.39	0.56
4:L:220:GLU:HG2	4:L:231:VAL:HG12	1.87	0.56
4:L:86:ARG:NE	4:L:107:ASP:OD2	2.36	0.55
6:R:176:SER:HA	6:R:198:THR:HB	1.88	0.55
4:L:214:HIS:O	4:L:236:ARG:NH1	2.36	0.55

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:H:51:PHE:HE1	5:H:119:ILE:HD12	1.72	0.54
6:R:195:HIS:O	6:R:200:ASN:ND2	2.41	0.54
4:L:90:SER:OG	4:L:97:THR:O	2.24	0.54
6:R:20:VAL:HG12	7:C:1:NAG:H82	1.89	0.54
5:H:110:THR:HG23	5:H:137:THR:HA	1.90	0.54
2:B:267:ASP:N	2:B:267:ASP:OD1	2.39	0.54
1:A:190:THR:HG22	1:A:195:HIS:HA	1.90	0.54
4:L:192:ASP:OD1	4:L:194:LYS:NZ	2.40	0.54
2:B:142:HIS:ND1	2:B:163:ASP:OD2	2.36	0.53
5:H:220:VAL:H	5:H:237:LEU:HD22	1.72	0.53
4:L:133:ARG:NH1	4:L:196:SER:O	2.42	0.53
1:A:271:LYS:HE2	1:A:323:PHE:HB3	1.91	0.53
1:A:339:VAL:HA	1:A:342:VAL:HG22	1.91	0.53
5:H:59:ALA:HB3	5:H:62:LYS:HB2	1.91	0.53
5:H:137:THR:HG21	5:H:173:PHE:HZ	1.74	0.52
2:B:145:TYR:O	2:B:162:GLY:N	2.42	0.52
1:A:207:GLU:HA	1:A:210:LYS:HD3	1.90	0.52
4:L:143:PHE:HB2	4:L:158:VAL:HG12	1.92	0.52
5:H:37:ARG:H	5:H:101:GLN:HE22	1.57	0.52
2:B:29:THR:OG1	2:B:30:LEU:N	2.43	0.52
5:H:146:PRO:HB3	5:H:172:TYR:HB3	1.91	0.52
2:B:228:ASP:OD1	2:B:228:ASP:N	2.43	0.51
4:L:88:SER:OG	4:L:99:THR:OG1	2.28	0.51
2:B:333:ASP:OD1	2:B:333:ASP:N	2.43	0.51
4:L:146:SER:OG	4:L:149:GLN:NE2	2.43	0.51
4:L:180:ARG:HH11	4:L:182:ASN:HB2	1.75	0.51
1:A:263:SER:HA	1:A:317:LYS:HD2	1.93	0.51
2:B:62:HIS:CD2	2:B:63:TRP:H	2.29	0.51
2:B:49:ARG:NH2	3:G:63:PRO:O	2.44	0.50
2:B:144:GLY:O	2:B:161:SER:OG	2.28	0.50
1:A:322:HIS:NE2	1:A:331:ASN:OD1	2.45	0.50
2:B:33:ILE:HG22	3:G:41:ARG:HH22	1.76	0.50
2:B:56:ALA:O	2:B:334:SER:OG	2.24	0.50
6:R:317:MET:HA	6:R:320:THR:HB	1.93	0.50
2:B:129:ARG:NH1	2:B:134:ARG:O	2.44	0.50
5:H:102:MET:HB3	5:H:105:LEU:HD21	1.94	0.50
4:L:92:SER:OG	4:L:95:ALA:N	2.42	0.50
6:R:157:VAL:O	6:R:160:THR:OG1	2.25	0.50
5:H:28:GLY:O	5:H:37:ARG:NH1	2.44	0.49
6:R:5:GLU:HA	6:R:10:TYR:HD1	1.77	0.49
2:B:103:CYS:HA	2:B:113:ALA:O	2.12	0.49

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:332:VAL:O	1:A:336:PHE:HB2	2.12	0.49
1:A:262:THR:O	1:A:317:LYS:NZ	2.39	0.49
5:H:164:THR:HA	5:H:209:THR:HA	1.95	0.49
6:R:178:TYR:HA	6:R:188:GLY:O	2.13	0.49
2:B:31:SER:O	2:B:34:THR:OG1	2.26	0.49
4:L:163:ASN:HA	4:L:197:THR:HB	1.94	0.49
2:B:37:ILE:O	2:B:301:LYS:NZ	2.37	0.49
2:B:294:CYS:HB3	2:B:308:LEU:HB2	1.95	0.49
5:H:170:LYS:NZ	5:H:198:GLN:OE1	2.46	0.49
6:R:183:MET:HG2	6:R:285:PRO:HB2	1.95	0.49
2:B:254:ASP:OD1	2:B:255:LEU:N	2.45	0.49
5:H:146:PRO:HD3	5:H:226:HIS:CD2	2.48	0.49
6:R:304:VAL:HA	6:R:308:MET:HB2	1.95	0.49
2:B:58:ILE:HG12	2:B:74:SER:HB3	1.95	0.48
2:B:237:ASN:OD1	2:B:238:GLY:N	2.47	0.48
4:L:154:GLY:CA	4:L:206:LEU:O	2.62	0.48
2:B:129:ARG:NH2	2:B:136:SER:OG	2.46	0.48
2:B:248:ALA:HA	2:B:273:ILE:HG13	1.95	0.48
4:L:141:SER:O	4:L:159:CYS:HA	2.13	0.48
2:B:153:ASP:OD1	2:B:153:ASP:N	2.43	0.48
6:R:181:GLU:H	6:R:184:GLN:HA	1.79	0.48
6:R:326:ASN:O	6:R:328:LEU:N	2.47	0.48
6:R:316:CYS:O	6:R:320:THR:OG1	2.26	0.48
5:H:112:MET:HA	5:H:134:THR:O	2.14	0.48
6:R:18:GLY:O	6:R:21:ARG:NH1	2.45	0.48
1:A:26:ASP:OD2	2:B:78:LYS:HE2	2.13	0.47
4:L:155:ALA:HB3	4:L:206:LEU:HB2	1.96	0.47
6:R:196:GLU:O	6:R:199:ASN:ND2	2.47	0.47
1:A:12:ALA:O	2:B:88:ASN:ND2	2.48	0.47
2:B:60:ALA:O	2:B:73:ALA:N	2.41	0.47
3:G:41:ARG:O	3:G:44:VAL:N	2.48	0.47
4:L:51:HIS:ND1	4:L:53:ASN:OD1	2.47	0.47
5:H:86:ARG:HD2	5:H:104:SER:HB2	1.95	0.47
5:H:215:TRP:CD1	5:H:216:PRO:HA	2.49	0.47
1:A:264:ILE:H	1:A:317:LYS:HD2	1.80	0.47
5:H:92:ASP:OD2	5:H:95:LYS:HD2	2.15	0.47
6:R:8:ASN:HD22	6:R:107:PRO:HG3	1.79	0.47
1:A:47:SER:HA	1:A:50:VAL:HG12	1.96	0.47
5:H:54:HIS:NE2	5:H:120:TYR:OH	2.38	0.47
1:A:250:PHE:HE1	1:A:264:ILE:HG13	1.79	0.46
4:L:26:GLN:OE1	4:L:126:GLY:N	2.47	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:147:SER:OG	2:B:160:SER:OG	2.28	0.46
1:A:38:LEU:HG	1:A:46:LYS:HD2	1.96	0.46
2:B:200:VAL:HG12	2:B:210:LEU:HA	1.96	0.46
2:B:81:ILE:HD12	2:B:90:VAL:HB	1.98	0.46
6:R:326:ASN:HD22	6:R:326:ASN:HA	1.54	0.46
1:A:25:GLU:HG2	1:A:28:GLU:OE2	2.16	0.46
4:L:221:ALA:O	4:L:229:PRO:HA	2.15	0.46
2:B:119:ASN:ND2	2:B:143:THR:O	2.45	0.46
1:A:311:ASN:O	1:A:313:ARG:NH1	2.49	0.46
2:B:227:SER:OG	2:B:228:ASP:OD1	2.33	0.46
4:L:215:ASN:OD1	4:L:237:ASN:N	2.48	0.46
2:B:148:CYS:HG	2:B:189:SER:HG	1.62	0.45
1:A:257:LYS:NZ	6:R:335:THR:H	2.14	0.45
6:R:163:MET:SD	6:R:163:MET:N	2.89	0.45
6:R:271:VAL:HG21	6:R:291:PRO:HG3	1.98	0.45
5:H:181:TRP:O	5:H:185:SER:N	2.47	0.45
6:R:311:LYS:HA	6:R:314:ARG:HG3	1.98	0.45
1:A:218:VAL:HG12	1:A:220:ALA:H	1.80	0.45
6:R:175:TRP:CD1	6:R:199:ASN:HB2	2.51	0.45
2:B:300:LEU:HD11	3:G:44:VAL:HG11	1.98	0.45
6:R:70:THR:OG1	6:R:73:ASN:ND2	2.50	0.45
5:H:51:PHE:CE2	5:H:117:ARG:HD2	2.51	0.45
1:A:18:MET:HA	1:A:21:ARG:HH21	1.82	0.45
2:B:80:ILE:HA	2:B:91:HIS:O	2.17	0.45
4:L:58:LEU:HB3	4:L:76:MET:HB2	1.99	0.45
4:L:71:LEU:HD22	5:H:128:ASP:HA	1.99	0.45
4:L:195:ASP:N	4:L:195:ASP:OD1	2.49	0.45
6:R:269:ALA:O	6:R:273:PHE:CB	2.65	0.45
4:L:170:ASN:HD22	4:L:171:VAL:H	1.64	0.44
2:B:189:SER:OG	2:B:190:LEU:N	2.49	0.44
4:L:133:ARG:HH12	4:L:197:THR:HG22	1.83	0.44
4:L:22:ILE:HD12	4:L:118:GLU:OE2	2.16	0.44
1:A:11:ALA:HA	1:A:14:GLU:HB3	1.99	0.44
1:A:209:LYS:HG2	1:A:213:HIS:ND1	2.32	0.44
1:A:351:CYS:O	6:R:135:ARG:NE	2.50	0.44
2:B:168:LEU:HD22	2:B:169:TRP:H	1.83	0.44
4:L:180:ARG:NH2	4:L:210:GLU:OE2	2.50	0.44
6:R:52:PHE:HD1	6:R:87:VAL:HG11	1.83	0.44
2:B:143:THR:OG1	2:B:144:GLY:N	2.51	0.44
2:B:209:LYS:HG2	2:B:221:THR:HA	2.00	0.44
4:L:143:PHE:HB3	5:H:151:LEU:HD12	1.99	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:H:174:PRO:HB3	5:H:228:ALA:HB2	1.99	0.44
6:R:181:GLU:N	6:R:184:GLN:HA	2.33	0.44
1:A:205:ARG:HE	1:A:245:GLU:HA	1.83	0.43
1:A:257:LYS:HG3	6:R:335:THR:HG22	2.00	0.43
2:B:70:LEU:O	2:B:81:ILE:HA	2.18	0.43
2:B:316:SER:OG	2:B:317:CYS:N	2.51	0.43
2:B:119:ASN:ND2	2:B:142:HIS:O	2.51	0.43
4:L:35:PRO:HG3	4:L:131:LEU:HD22	2.00	0.43
4:L:139:THR:O	4:L:161:LEU:HA	2.19	0.43
1:A:194:LEU:HD22	1:A:194:LEU:HA	1.92	0.43
5:H:182:ASN:OD1	5:H:221:THR:N	2.52	0.43
6:R:22:SER:HB3	6:R:25:GLU:HG2	2.00	0.43
2:B:79:LEU:O	2:B:92:ALA:HA	2.18	0.43
2:B:237:ASN:OD1	2:B:239:ASN:N	2.50	0.43
6:R:195:HIS:NE2	6:R:197:GLU:HB2	2.33	0.43
1:A:305:CYS:O	1:A:309:ASP:N	2.51	0.43
1:A:322:HIS:HB2	1:A:334:PHE:CE2	2.54	0.43
2:B:81:ILE:HD11	2:B:91:HIS:CD2	2.54	0.43
6:R:75:ILE:HD11	6:R:127:SER:HA	2.00	0.43
6:R:109:GLY:HA2	6:R:112:LEU:HD12	1.98	0.43
1:A:38:LEU:HB3	1:A:46:LYS:HZ2	1.83	0.43
2:B:340:ASN:OD1	3:G:62:ASN:ND2	2.52	0.43
3:G:32:LEU:HD23	3:G:33:VAL:H	1.84	0.43
2:B:108:SER:OG	2:B:109:GLY:N	2.52	0.42
4:L:61:PHE:HB2	4:L:112:TYR:HB2	2.00	0.42
6:R:130:VAL:HG12	6:R:133:ILE:HD12	2.00	0.42
2:B:66:ASP:OD1	2:B:66:ASP:N	2.38	0.42
1:A:29:LYS:HD3	1:A:32:ARG:HD3	2.00	0.42
2:B:135:VAL:HG22	2:B:137:ARG:H	1.84	0.42
2:B:205:ASP:N	2:B:205:ASP:OD1	2.44	0.42
2:B:318:LEU:H	2:B:318:LEU:HG	1.77	0.42
5:H:220:VAL:HB	5:H:237:LEU:HD13	2.01	0.42
5:H:55:TRP:CD2	5:H:100:LEU:HD22	2.55	0.42
1:A:48:THR:HG21	1:A:327:THR:HG23	2.02	0.42
1:A:230:TYR:HD2	1:A:277:LYS:HE3	1.85	0.42
5:H:153:PRO:HB3	5:H:157:ASP:OD2	2.19	0.42
6:R:189:ILE:HD11	8:R:401:RET:H183	2.01	0.42
6:R:290:ILE:H	6:R:290:ILE:HG13	1.64	0.42
1:A:314:LYS:HG2	1:A:315:ASP:H	1.84	0.42
2:B:255:LEU:HB3	2:B:256:ARG:HD3	2.02	0.42
1:A:306:GLN:NE2	1:A:309:ASP:OD2	2.52	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:106:ALA:HB2	2:B:111:TYR:HB2	2.01	0.42
1:A:46:LYS:HA	1:A:49:ILE:HD12	2.02	0.41
2:B:63:TRP:HE1	2:B:319:GLY:C	2.22	0.41
6:R:195:HIS:CD2	6:R:198:THR:H	2.38	0.41
6:R:195:HIS:HB3	6:R:200:ASN:HD21	1.85	0.41
2:B:246:ASP:HA	2:B:273:ILE:H	1.86	0.41
2:B:239:ASN:HA	2:B:255:LEU:HD22	2.02	0.41
6:R:119:LEU:HD13	6:R:168:ALA:HB2	2.02	0.41
2:B:160:SER:OG	2:B:160:SER:O	2.37	0.41
4:L:113:CYS:O	4:L:124:GLY:N	2.53	0.41
4:L:165:TYR:HA	4:L:166:PRO:HA	1.91	0.41
5:H:121:TYR:H	5:H:124:SER:HG	1.67	0.41
5:H:174:PRO:HD3	5:H:226:HIS:NE2	2.36	0.41
5:H:107:SER:HA	5:H:138:VAL:HB	2.01	0.41
6:R:262:LEU:HD13	6:R:262:LEU:HA	1.93	0.41
6:R:103:PHE:CZ	6:R:105:PHE:HB2	2.56	0.41
2:B:62:HIS:O	2:B:71:VAL:N	2.53	0.40
4:L:103:LEU:HD12	4:L:103:LEU:HA	1.91	0.40
4:L:147:SER:HA	4:L:150:LEU:HD12	2.03	0.40
2:B:273:ILE:HG12	2:B:289:TYR:CD1	2.56	0.40
3:G:38:GLU:HA	3:G:41:ARG:HH11	1.86	0.40
1:A:208:ARG:O	1:A:212:ILE:HG22	2.22	0.40
1:A:16:SER:HA	1:A:19:ILE:HG22	2.03	0.40
2:B:190:LEU:HA	2:B:190:LEU:HD13	1.94	0.40
4:L:167:LYS:HB2	4:L:188:TRP:HD1	1.86	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	206/376 (55%)	169 (82%)	37 (18%)	0	100 100

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	B	310/340 (91%)	271 (87%)	39 (13%)	0	100	100
3	G	34/74 (46%)	29 (85%)	5 (15%)	0	100	100
4	L	215/239 (90%)	196 (91%)	19 (9%)	0	100	100
5	H	221/249 (89%)	201 (91%)	20 (9%)	0	100	100
6	R	332/348 (95%)	299 (90%)	33 (10%)	0	100	100
All	All	1318/1626 (81%)	1165 (88%)	153 (12%)	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	188/325 (58%)	185 (98%)	3 (2%)	62	79
2	B	260/283 (92%)	251 (96%)	9 (4%)	36	60
3	G	34/70 (49%)	34 (100%)	0	100	100
4	L	193/210 (92%)	190 (98%)	3 (2%)	62	79
5	H	191/215 (89%)	187 (98%)	4 (2%)	53	72
6	R	284/296 (96%)	280 (99%)	4 (1%)	67	81
All	All	1150/1399 (82%)	1127 (98%)	23 (2%)	61	74

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	184	ILE
1	A	194	LEU
1	A	313	ARG
2	B	47	THR
2	B	61	MET
2	B	82	TRP
2	B	96	ARG
2	B	134	ARG

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
2	B	200	VAL
2	B	256	ARG
2	B	266	HIS
2	B	340	ASN
4	L	133	ARG
4	L	161	LEU
4	L	170	ASN
5	H	176[A]	SER
5	H	176[B]	SER
5	H	223[A]	SER
5	H	223[B]	SER
6	R	147	ARG
6	R	163	MET
6	R	314	ARG
6	R	326	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	269	ASN
1	A	306	GLN
1	A	346	ASN
2	B	62	HIS
2	B	91	HIS
2	B	176	GLN
2	B	183	HIS
2	B	239	ASN
2	B	311	HIS
4	L	116	HIS
4	L	149	GLN
4	L	170	ASN
4	L	235	ASN
5	H	101	GLN
6	R	73	ASN
6	R	195	HIS
6	R	199	ASN
6	R	302	ASN
6	R	326	ASN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains (i)

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

5.5 Carbohydrates (i)

2 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	C	1	7,6	14,14,15	0.31	0	17,19,21	0.50	0
7	NAG	C	2	7	14,14,15	0.47	0	17,19,21	0.63	1 (5%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
7	NAG	C	1	7,6	-	0/6/23/26	0/1/1/1
7	NAG	C	2	7	-	2/6/23/26	0/1/1/1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
7	C	2	NAG	C1-O5-C5	2.14	115.10	112.19

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	C	2	NAG	O5-C5-C6-O6
7	C	2	NAG	C4-C5-C6-O6

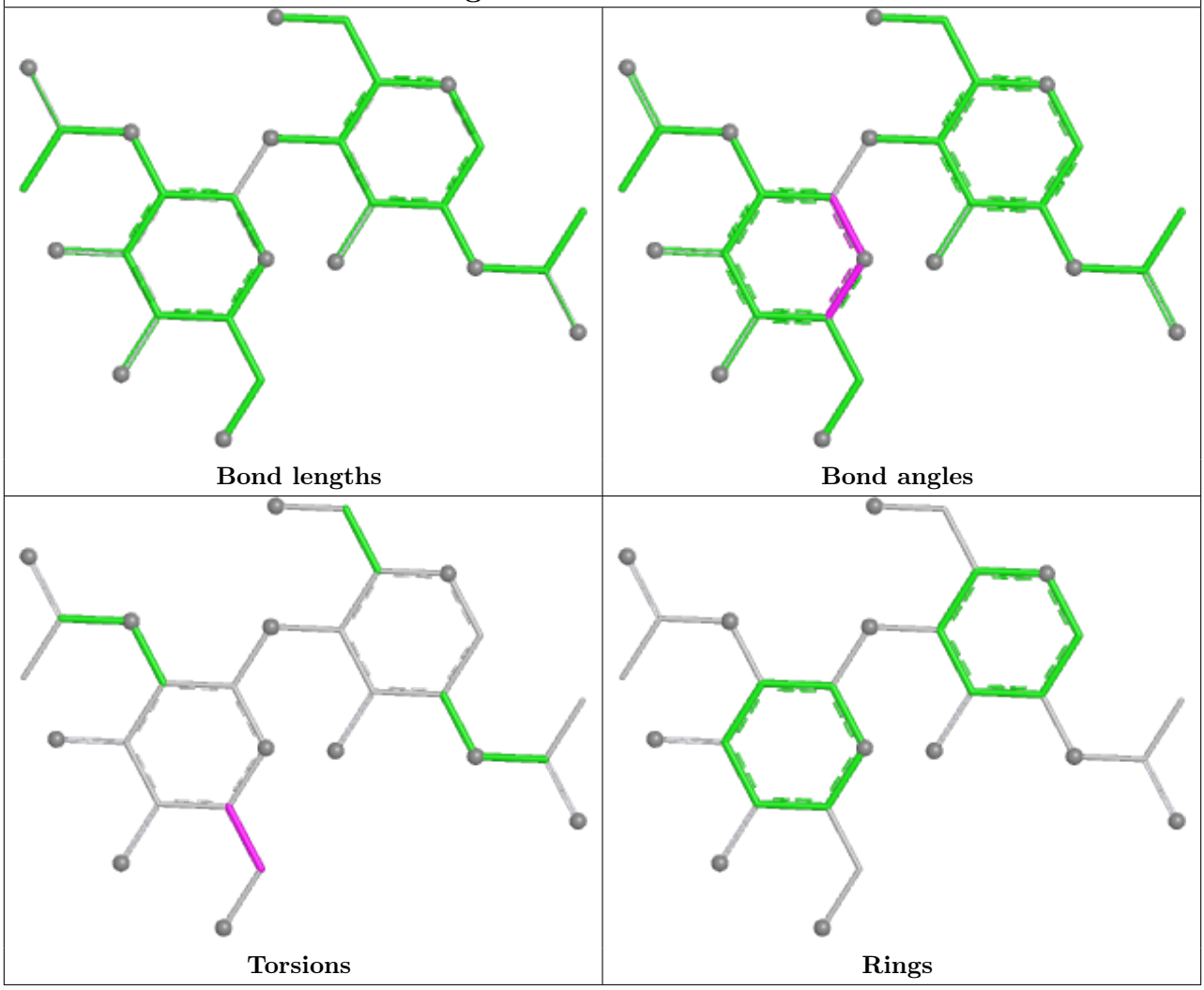
There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	C	1	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 C



5.6 Ligand geometry (i)

1 ligand is 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	$\#\|Z\| > 2$
8	RET	R	401	6	20,20,21	3.31	10 (50%)	27,27,28	1.61	7 (25%)

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	RET	R	401	6	-	5/13/30/31	0/1/1/1

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
8	R	401	RET	C8-C9	6.41	1.59	1.45
8	R	401	RET	C7-C6	5.94	1.66	1.45
8	R	401	RET	C1-C6	5.76	1.61	1.53
8	R	401	RET	C14-C13	4.30	1.37	1.33
8	R	401	RET	C18-C5	4.29	1.58	1.50
8	R	401	RET	C4-C5	4.18	1.59	1.51
8	R	401	RET	C11-C10	3.80	1.55	1.43
8	R	401	RET	C12-C13	3.69	1.53	1.45
8	R	401	RET	C2-C1	2.24	1.59	1.54
8	R	401	RET	C16-C1	2.10	1.57	1.53

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
8	R	401	RET	C19-C9-C10	-3.55	117.95	122.92
8	R	401	RET	C8-C9-C10	3.43	124.20	118.94
8	R	401	RET	C16-C1-C6	-2.87	105.64	110.30
8	R	401	RET	C17-C1-C6	2.75	114.76	110.30
8	R	401	RET	C18-C5-C4	-2.53	108.76	113.62
8	R	401	RET	C7-C8-C9	2.46	129.95	126.23
8	R	401	RET	C7-C6-C5	2.11	126.58	121.46

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
8	R	401	RET	C7-C8-C9-C10
8	R	401	RET	C7-C8-C9-C19
8	R	401	RET	C11-C12-C13-C14
8	R	401	RET	C11-C12-C13-C20
8	R	401	RET	C9-C10-C11-C12

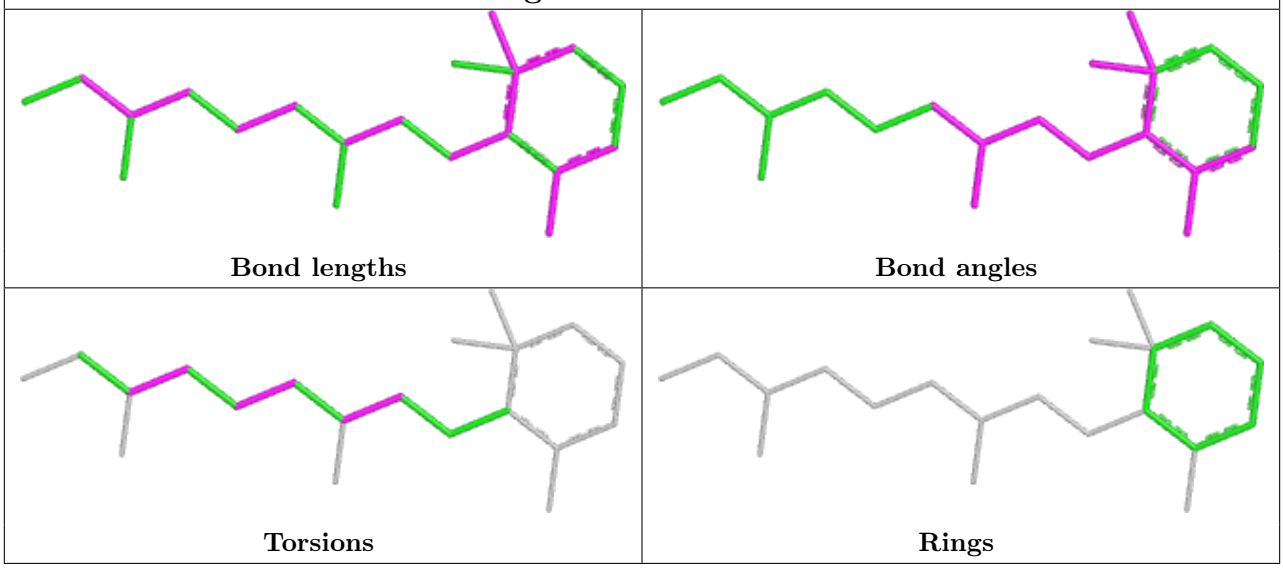
There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
8	R	401	RET	1	0

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

Ligand RET R 401



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.

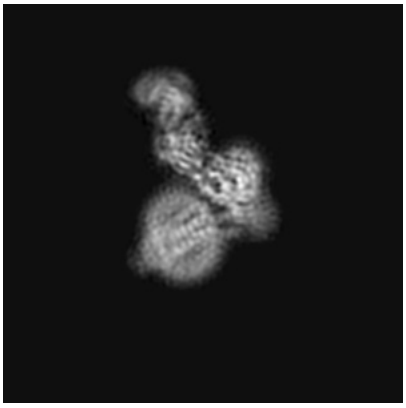
6 Map visualisation (i)

This section contains visualisations of the EMDB entry EMD-4598. These allow visual inspection of the internal detail of the map and identification of artifacts.

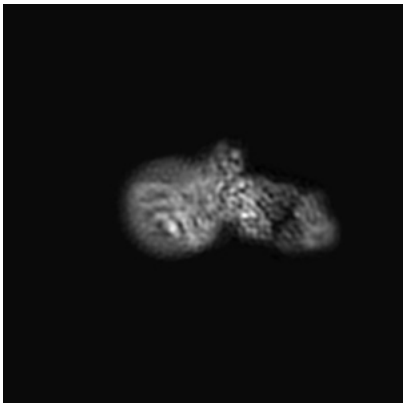
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections (i)

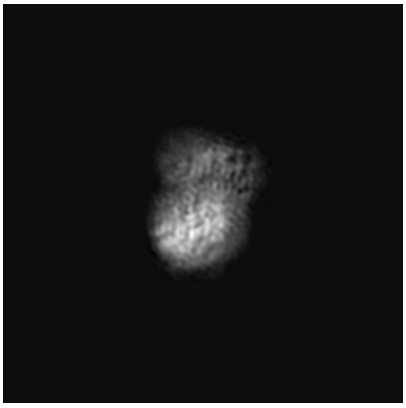
6.1.1 Primary map



X



Y

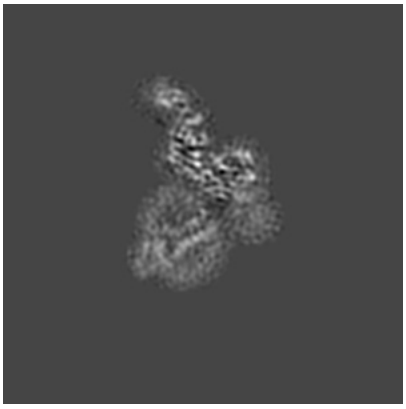


Z

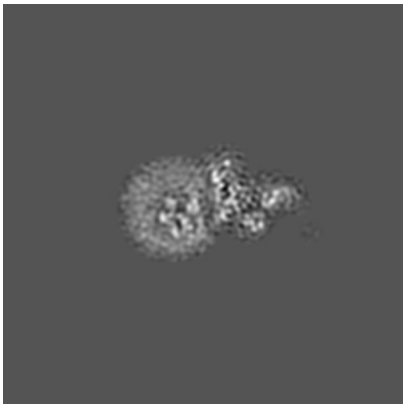
The images above show the map projected in three orthogonal directions.

6.2 Central slices (i)

6.2.1 Primary map



X Index: 169



Y Index: 169

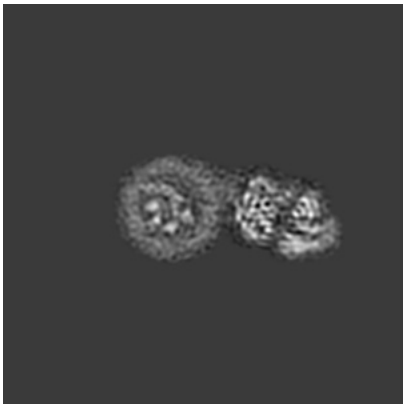
Z Index: 169

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

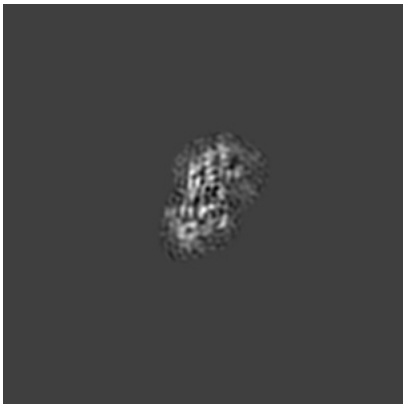
6.3 Largest variance slices (i)

6.3.1 Primary map

X Index: 159



Y Index: 146

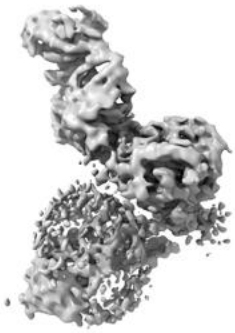


Z Index: 203

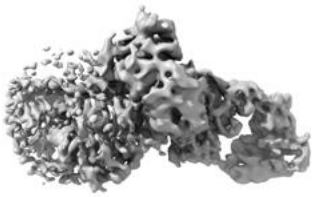
The images above show the largest variance slices of the map in three orthogonal directions.

6.4 Orthogonal surface views (i)

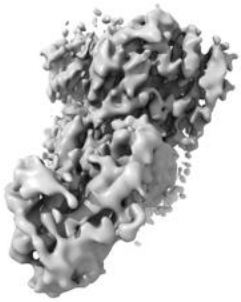
6.4.1 Primary map



X



Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.03. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

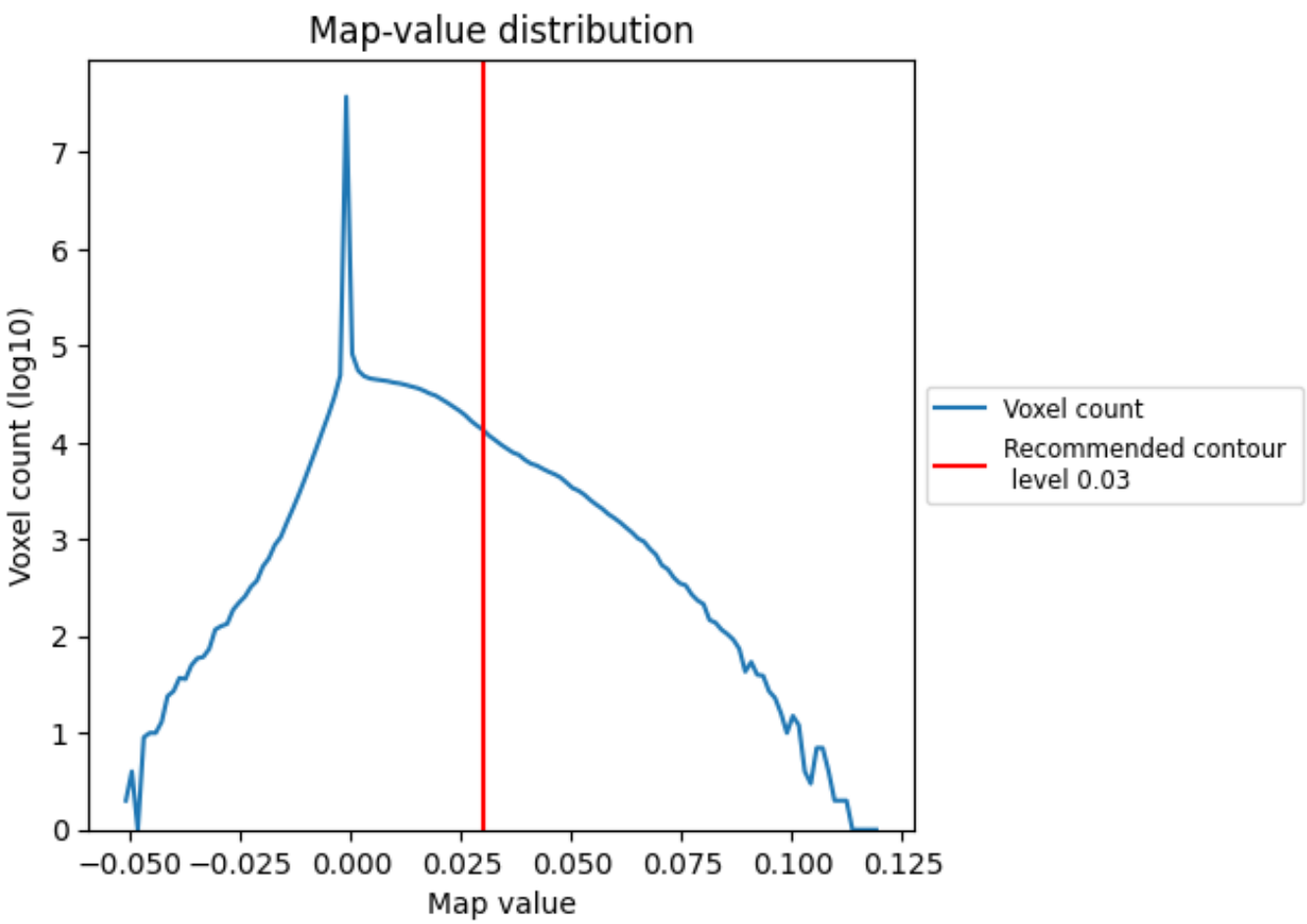
6.5 Mask visualisation (i)

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

7 Map analysis (i)

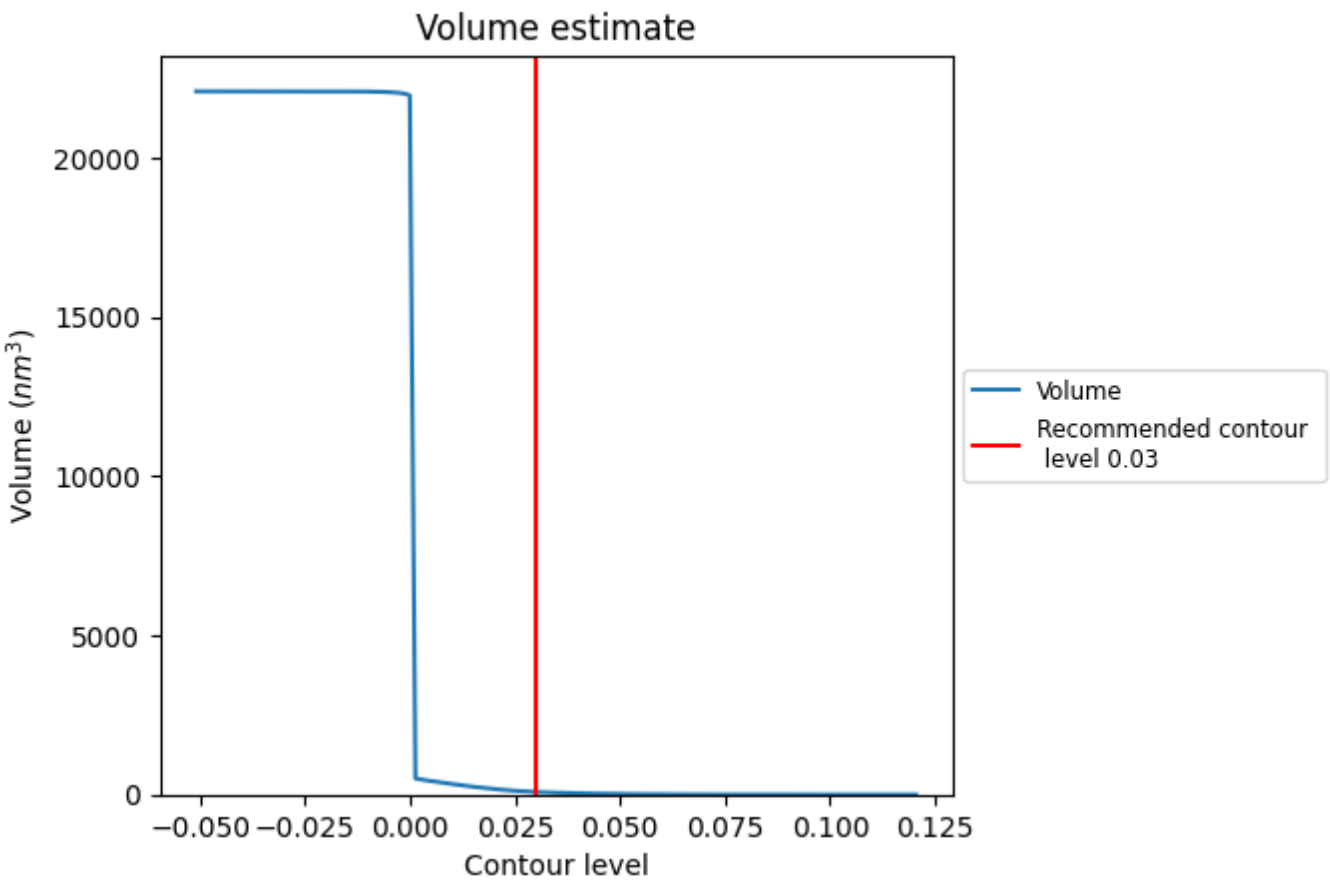
This section contains the results of statistical analysis of the map.

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.

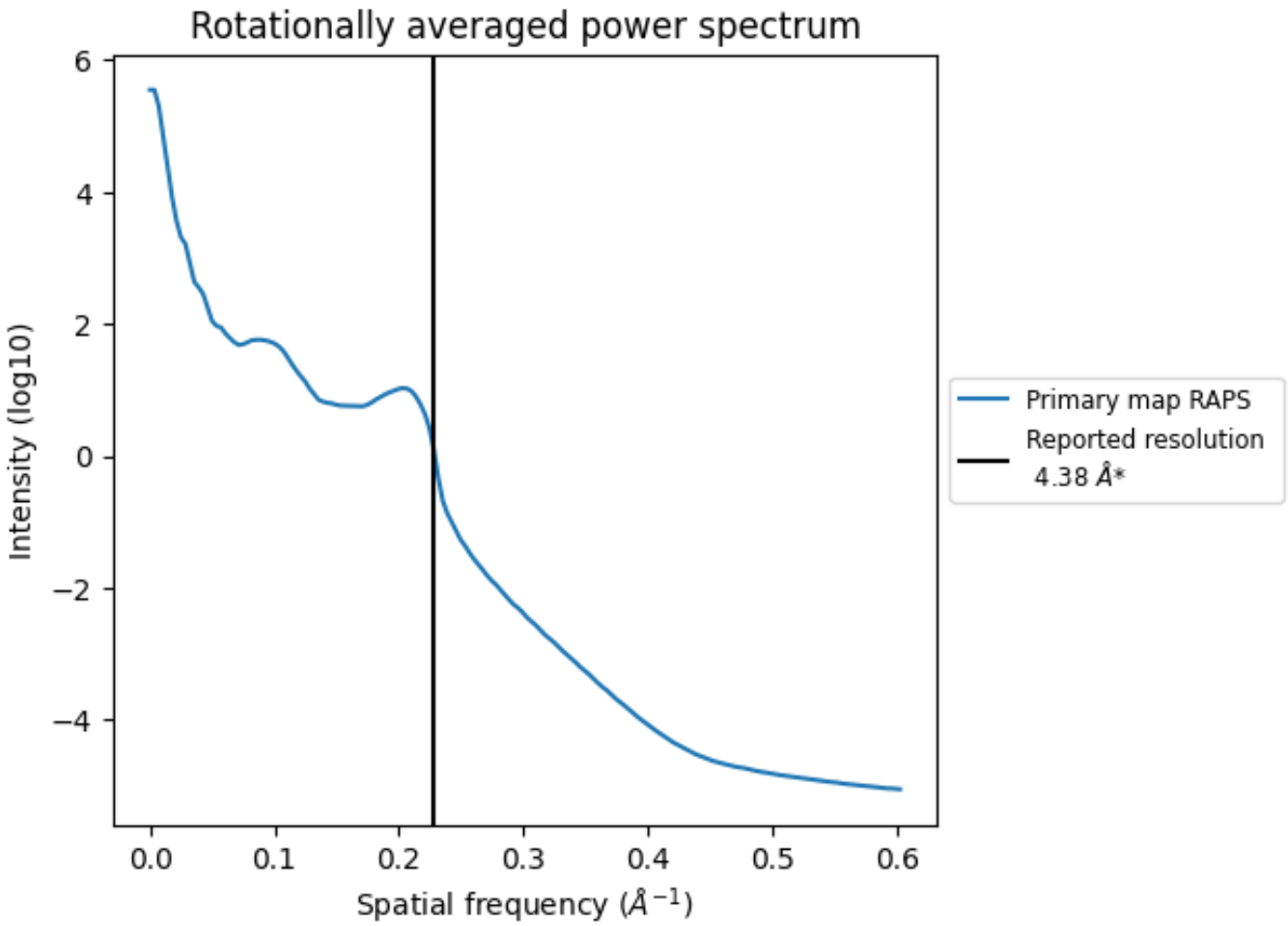
7.2 Volume estimate (i)



The volume at the recommended contour level is 82 nm^3; this corresponds to an approximate mass of 74 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 0.228 Å^{-1}

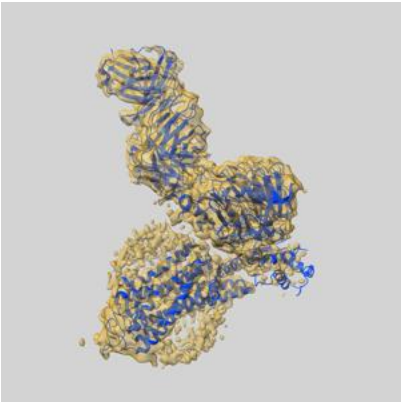
8 Fourier-Shell correlation (i)

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

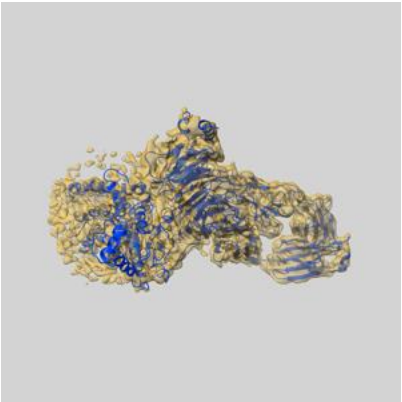
9 Map-model fit (i)

This section contains information regarding the fit between EMDB map EMD-4598 and PDB model 6QNO. Per-residue inclusion information can be found in section 3 on page 6.

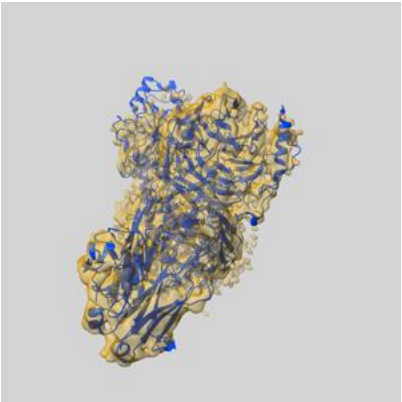
9.1 Map-model overlay (i)



X



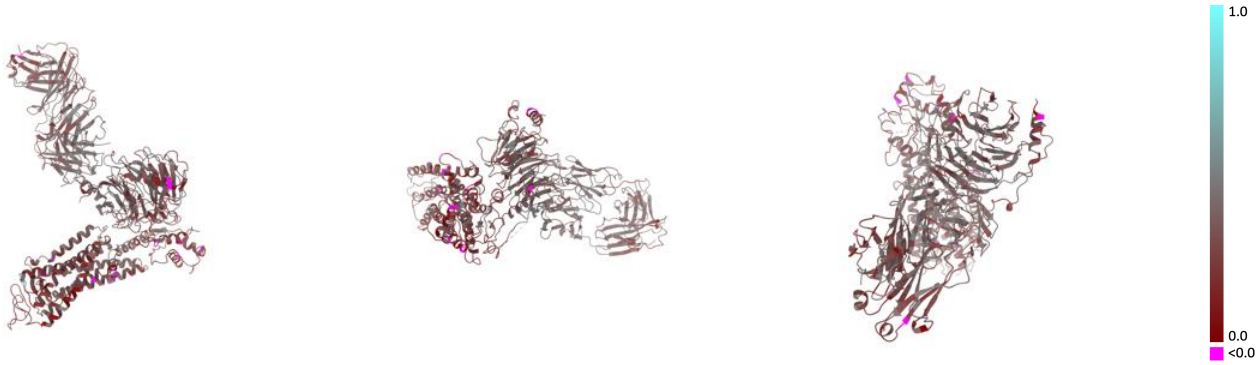
Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.03 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.

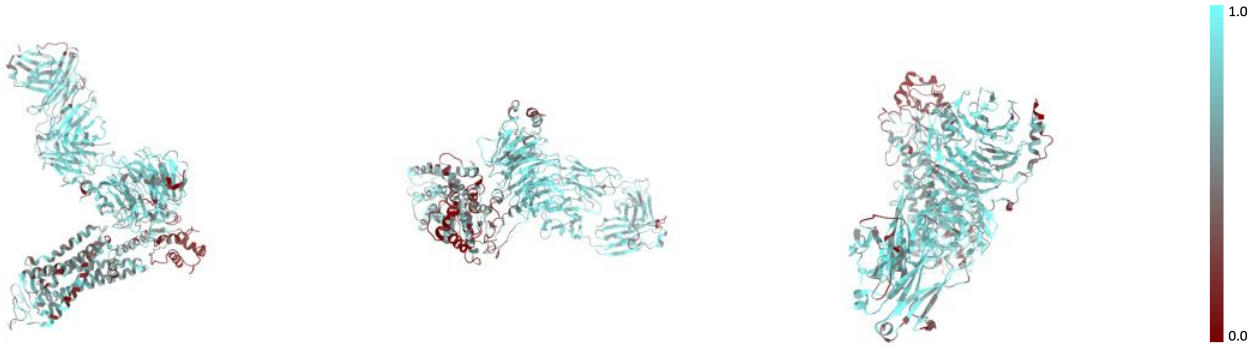
9.2 Q-score mapped to coordinate model (i)



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.

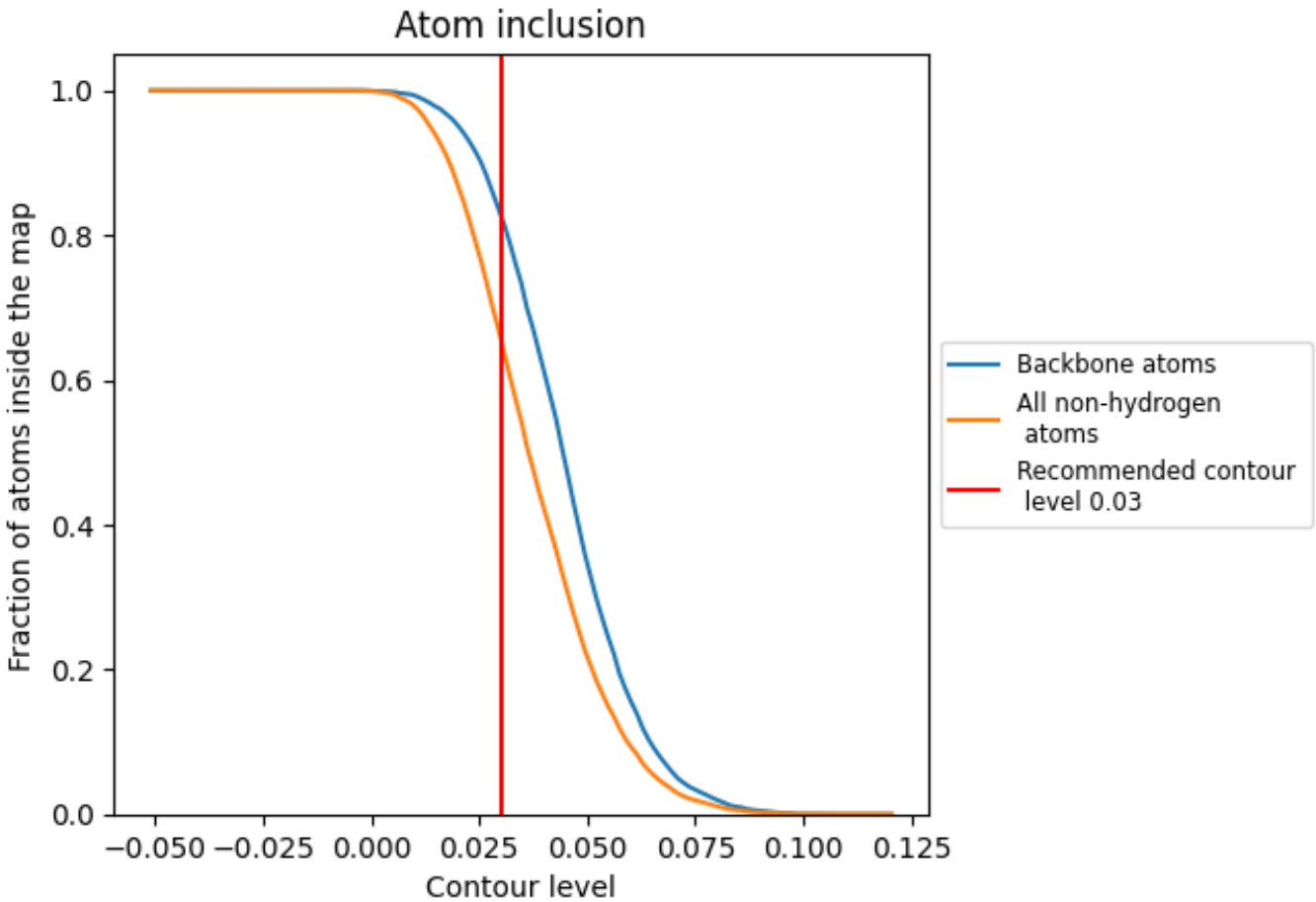
9.3 Atom inclusion mapped to coordinate model (i)



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 (0.03).

9.4 Atom inclusion (i)



At the recommended contour level, 83% of all backbone atoms, 66% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary (i)

The table lists the average atom inclusion at the recommended contour level (0.03) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.6570	0.3230
A	0.3806	0.2640
B	0.7844	0.3680
C	0.2500	0.3450
G	0.5208	0.2870
H	0.7509	0.3560
L	0.7591	0.3620
R	0.6183	0.2800