

# Full wwPDB EM Validation Report (i)

#### Oct 8, 2022 – 06:00 PM EDT

PDB ID : 7T9M

EMDB ID : EMD-25762

Title: Human Thyrotropin receptor bound by CS-17 Inverse Agonist Fab/Org

274179-0 Antagonist

Authors: Faust, B.; Cheng, Y.; Manglik, A.

Deposited on : 2021-12-19

Resolution : 3.10 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 (i) symbol.

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

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

EMDB validation analysis : 0.0.1.dev43

Mogul : 1.8.5 (274361), CSD as541be (2020)

MolProbity: 4.02b-467

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)

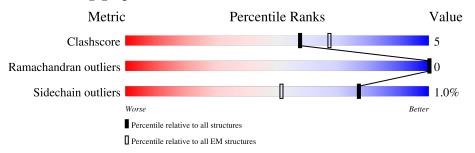
Validation Pipeline (wwPDB-VP) : 2.31.2

# 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 3.10 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 $(\#  ext{Entries})$	${ m EM~structures} \ (\#{ m Entries})$		
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	Н	217	8%	11%					
2	L	213	7% 85%	13% •					
3	R	702	74% 6%	20%					
4	A	2	100%						



# 2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 15615 atoms, of which 7788 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 CS-17 Heavy Chain.

Mol	Chain	Residues	Atoms						AltConf	Trace
1	Н	217	Total 3219	C 1035	H 1583	N 260	O 334	S 7	0	0

• Molecule 2 is a protein called CS-17 Light Chain.

Mol	Chain	Residues	Atoms						AltConf	Trace
9	ī	213	Total	С	Н	N	О	S	0	0
	ь	213	3186	1017	1554	272	333	10	0	

• Molecule 3 is a protein called Thyrotropin receptor.

Mol	Chain	Residues	Atoms						AltConf	Trace
3	R	565	Total 9071	C 2947	H 4582	N 718	O 791	S 33	0	0

There are 59 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
R	13	ASP	-	expression tag	UNP A0A0A0MTJ0
R	14	TYR	-	expression tag	UNP A0A0A0MTJ0
R	15	LYS	-	expression tag	UNP A0A0A0MTJ0
R	16	ASP	-	expression tag	UNP A0A0A0MTJ0
R	17	ASP	-	expression tag	UNP A0A0A0MTJ0
R	18	ASP	-	expression tag	UNP A0A0A0MTJ0
R	19	ASP	-	expression tag	UNP A0A0A0MTJ0
R	20	GLY	-	expression tag	UNP A0A0A0MTJ0
R	21	THR	-	expression tag	UNP A0A0A0MTJ0
R	?	-	ALA	deletion	UNP A0A0A0MTJ0
R	?	-	LEU	deletion	UNP A0A0A0MTJ0
R	?	-	ASN	deletion	UNP A0A0A0MTJ0
R	?	-	SER	deletion	UNP A0A0A0MTJ0
R	?	-	PRO	deletion	UNP A0A0A0MTJ0
R	?	-	LEU	deletion	UNP A0A0A0MTJ0



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Chain	Residue	Modelled	Actual	Comment	Reference
R	?	-	HIS	deletion	UNP A0A0A0MTJ0
R	?	-	GLN	deletion	UNP A0A0A0MTJ0
R	?	-	GLU	deletion	UNP A0A0A0MTJ0
R	?	-	TYR	deletion	UNP A0A0A0MTJ0
R	?	-	GLU	deletion	UNP A0A0A0MTJ0
R	?	-	GLU	deletion	UNP A0A0A0MTJ0
R	?	-	ASN	deletion	UNP A0A0A0MTJ0
R	?	-	LEU	deletion	UNP A0A0A0MTJ0
R	?	-	GLY	deletion	UNP A0A0A0MTJ0
R	?	-	ASP	deletion	UNP A0A0A0MTJ0
R	?	-	SER	deletion	UNP A0A0A0MTJ0
R	?	-	ILE	deletion	UNP A0A0A0MTJ0
R	?	-	VAL	deletion	UNP A0A0A0MTJ0
R	?	-	GLY	deletion	UNP A0A0A0MTJ0
R	?	-	TYR	deletion	UNP A0A0A0MTJ0
R	?	-	LYS	deletion	UNP A0A0A0MTJ0
R	?	-	GLU	deletion	UNP A0A0A0MTJ0
R	?	-	LYS	deletion	UNP A0A0A0MTJ0
R	?	-	SER	deletion	UNP A0A0A0MTJ0
R	?	-	LYS	deletion	UNP A0A0A0MTJ0
R	?	-	PHE	deletion	UNP A0A0A0MTJ0
R	?	-	GLN	deletion	UNP A0A0A0MTJ0
R	?	-	ASP	deletion	UNP A0A0A0MTJ0
R	?	-	THR	deletion	UNP A0A0A0MTJ0
R	?	-	HIS	deletion	UNP A0A0A0MTJ0
R	?	-	ASN	deletion	UNP A0A0A0MTJ0
R	?	-	ASN	deletion	UNP A0A0A0MTJ0
R	?	-	ALA	deletion	UNP A0A0A0MTJ0
R	?	-	HIS	deletion	UNP A0A0A0MTJ0
R	?	-	TYR	deletion	UNP A0A0A0MTJ0
R	?	-	TYR	deletion	UNP A0A0A0MTJ0
R	?	-	VAL	deletion	UNP A0A0A0MTJ0
R	?	-	PHE	deletion	UNP A0A0A0MTJ0
R	?	-	PHE	deletion	UNP A0A0A0MTJ0
R	?	-	GLU	deletion	UNP A0A0A0MTJ0
R	?	-	GLU	deletion	UNP A0A0A0MTJ0
R	?	-	GLN	deletion	UNP A0A0A0MTJ0
R	?	-	GLU	deletion	UNP A0A0A0MTJ0
R	?	-	ASP	deletion	UNP A0A0A0MTJ0
R	?	-	GLU	deletion	UNP A0A0A0MTJ0
R	?	-	ILE	deletion	UNP A0A0A0MTJ0
R	?	_	ILE	deletion	UNP A0A0A0MTJ0



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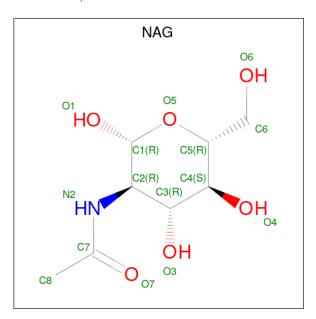
Chain	Residue	Modelled	Actual	Comment	Reference
R	?	-	GLY	deletion	UNP A0A0A0MTJ0
R	?	-	PHE	deletion	UNP A0A0A0MTJ0

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



Mol	Chain	Residues	Atoms					AltConf	Trace
4	A	2	Total 55		H 27	N 2	O 10	0	0

 $\bullet$  Molecule 5 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula:  $\rm C_8H_{15}NO_6).$ 



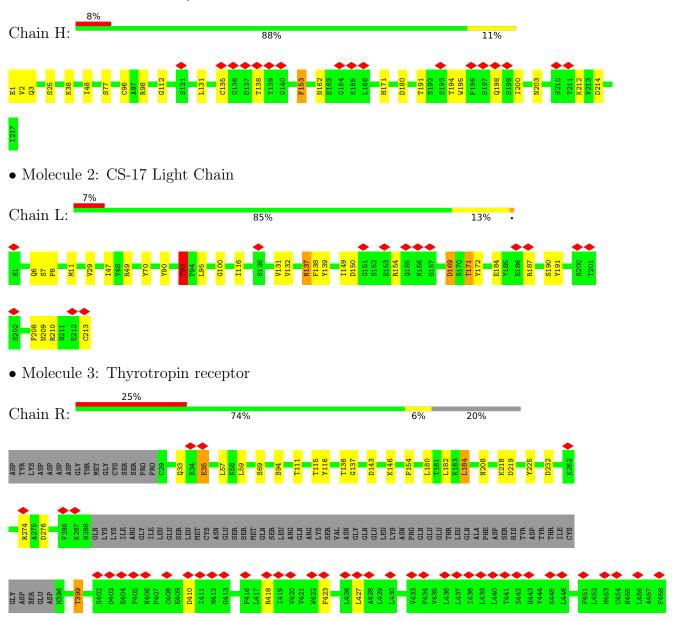
Mol	Chain	Residues	Atoms					AltConf
5	R	1	Total	С	Н	N	О	0
	16	1	84	24	42	3	15	U
5	R	1	Total	С	Н	N	О	0
9	10	1	84	24	42	3	15	U
5	R.	1	Total	С	Н	N	О	0
5	n	1	84	24	42	3	15	U



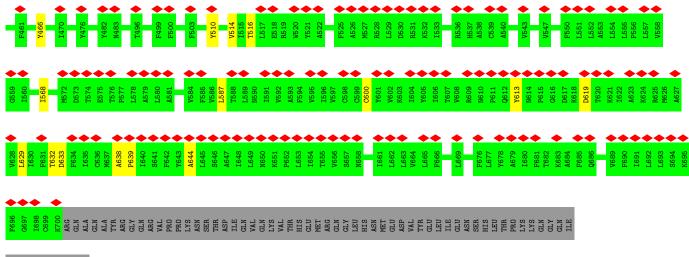
# 3 Residue-property plots (i)

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: CS-17 Heavy Chain







SER GLU GLU TYR MET GLN THR VAL LEU

 $\bullet$  Molecule 4: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain A: 100%





# 4 Experimental information (i)

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	41054	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE	Depositor
	CORRECTION	
Microscope	TFS KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose $(e^-/\text{Å}^2)$	77	Depositor
Minimum defocus (nm)	800	Depositor
Maximum defocus (nm)	2200	Depositor
Magnification	130000	Depositor
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	0.934	Depositor
Minimum map value	-0.333	Depositor
Average map value	-0.001	Depositor
Map value standard deviation	0.023	Depositor
Recommended contour level	0.2	Depositor
Map size (Å)	288.512, 288.512, 288.512	wwPDB
Map dimensions	448, 448, 448	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.644, 0.644, 0.644	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

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		
IVIOI	Chain	RMSZ	# Z  > 5	RMSZ	# Z  > 5	
1	Н	0.31	0/1677	0.55	1/2292~(0.0%)	
2	L	0.38	0/1673	0.64	1/2274~(0.0%)	
3	R	0.32	0/4600	0.54	1/6263 (0.0%)	
All	All	0.33	0/7950	0.56	3/10829 (0.0%)	

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	$\mathbf{Z}$	$Observed(^o)$	$\mathrm{Ideal}(^{o})$
1	Н	153	PHE	C-N-CD	-6.01	107.37	120.60
3	R	399	THR	N-CA-C	5.32	125.36	111.00
2	L	93	TYR	N-CA-C	5.19	125.01	111.00

There are no chirality outliers.

There are no planarity outliers.

#### 5.2 Too-close contacts (i)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	Н	1636	1583	1583	15	0
2	L	1632	1554	1555	31	0
3	R	4489	4582	4576	35	0
4	A	28	27	25	0	0
5	R	42	42	39	0	0



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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	7827	7788	7778	78	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 (78) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

2:L:6:GLN:OE1 2:L:10 3:R:466:TYR:OH 3:R: 2:L:6:GLN:OE1 2:L:1 2:L:150:ASP:OD1 2:L:19	100:GLY:CA 568:ILE:O 100:GLY:N 90:SER:OG 9:ASP:OD2	2.37 2.08 2.25 2.11	0.73 0.71 0.69
3:R:466:TYR:OH 3:R: 2:L:6:GLN:OE1 2:L:1 2:L:150:ASP:OD1 2:L:1	568:ILE:O 100:GLY:N 90:SER:OG	2.08 2.25	0.71 0.69
2:L:6:GLN:OE1 2:L:1 2:L:150:ASP:OD1 2:L:19	100:GLY:N 90:SER:OG	2.25	0.69
2:L:150:ASP:OD1 2:L:19	90:SER:OG		
		2.11	
2.D.612.TVD.ОЦ 2.D.61	9:ASP:OD2		0.67
9.0.019.11.0.01   9.0.01		2.14	0.63
3:R:33:GLN:HG3 3:R:	33:GLN:O	1.99	0.63
3:R:184:LEU:HD22 3:R:1	184:LEU:H	1.63	0.62
2:L:184:GLU:OE2 2:L:18	7:ARG:NH1	2.31	0.62
3:R:115:THR:HG22 3:R:1	137:GLY:O	2.00	0.61
2:L:137:ASN:N 2:L:137	7:ASN:HD22	1.97	0.61
3:R:276:ASP:OD1 3:R:39	9:THR:OG1	2.20	0.60
2:L:29:VAL:O 2:L:7	0:TYR:OH	2.21	0.59
3:R:182:LEU:HB3 3:R:184	4:LEU:HD21	1.85	0.59
2:L:8:PRO:CG 2:L:11	1:MET:HE2	2.32	0.59
3:R:184:LEU:HD22 3:R:1	184:LEU:N	2.18	0.59
2:L:47:ILE:HG22 2:L:4	49:ARG:O	2.04	0.57
1:H:153:PHE:CE2 1:H:1	180:ASP:O	2.58	0.57
2:L:8:PRO:HG3 2:L:11	1:MET:HE2	1.87	0.56
2:L:138:PHE:O 2:L:171	1:THR:HG22	2.06	0.56
1:H:198:GLN:OE1 1:H:20	0:ILE:HD13	2.06	0.55
1:H:138:THR:HG21 1:H:19	95:TRP:HZ3	1.72	0.55
2:L:139:TYR:O 2:L:17	2:TYR:HD2	1.90	0.54
2:L:6:GLN:OE1 2:L:1	100:GLY:C	2.47	0.53
2:L:93:TYR:O 2:L:93	3:TYR:HD1	1.91	0.53
2:L:138:PHE:C 2:L:171	I:THR:HG22	2.29	0.53
2:L:191:TYR:N 2:L:2	208:PHE:O	2.41	0.53
3:R:35:GLU:N 3:R:35	5:GLU:OE2	2.41	0.52
3:R:276:ASP:HA 3:R:39	9:THR:OG1	2.09	0.52
1:H:153:PHE:CD2 1:H:1	180:ASP:O	2.63	0.52
3:R:57:LEU:HG 3:R:59	0:LEU:HD21	1.91	0.52
3:R:276:ASP:HA 3:R:39	99:THR:CB	2.38	0.52
1:H:131:LEU:HD13 2:L:132	2:VAL:HG21	1.92	0.52
1:H:2:VAL:HG11 1:H:98	8:ARG:NH1	2.25	0.51



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Continued from prev		Interatomic	Clash
Atom-1	Atom-2	${\rm distance}\ (\mathring{\rm A})$	overlap(Å)
1:H:212:LYS:NZ	1:H:214:ASP:OD1	2.43	0.51
3:R:276:ASP:HA	3:R:399:THR:HB	1.92	0.50
3:R:629:LEU:O	3:R:632:THR:OG1	2.25	0.50
3:R:154:PHE:HB3	3:R:180:LEU:HD12	1.92	0.50
3:R:69:SER:OG	3:R:94:SER:O	2.16	0.49
3:R:423:PHE:CZ	3:R:427:LEU:HD11	2.47	0.49
3:R:208:ASN:ND2	3:R:232:ASP:OD2	2.45	0.49
1:H:162:ASN:OD1	1:H:203:ASN:ND2	2.45	0.48
3:R:629:LEU:HD23	3:R:633:ASP:OD2	2.13	0.48
2:L:8:PRO:CG	2:L:11:MET:CE	2.91	0.48
1:H:38:LYS:HB2	1:H:48:ILE:HD11	1.96	0.48
2:L:116:ILE:HD11	2:L:131:VAL:CG1	2.44	0.47
2:L:139:TYR:O	2:L:172:TYR:CD2	2.67	0.47
3:R:143:ASP:OD2	3:R:146:LYS:NZ	2.48	0.47
3:R:57:LEU:CG	3:R:59:LEU:HD21	2.45	0.46
2:L:169:ASP:N	2:L:169:ASP:OD1	2.48	0.46
3:R:116:TYR:O	3:R:116:TYR:CD2	2.69	0.46
3:R:225:TYR:C	3:R:225:TYR:CD1	2.89	0.46
2:L:137:ASN:N	2:L:137:ASN:ND2	2.62	0.45
2:L:139:TYR:O	2:L:139:TYR:CD1	2.69	0.45
3:R:510:TYR:O	3:R:514:VAL:HG23	2.17	0.45
3:R:225:TYR:CD1	3:R:225:TYR:O	2.69	0.45
3:R:638:ALA:HB3	3:R:639:PRO:HD3	1.98	0.45
1:H:191:THR:O	1:H:194:THR:OG1	2.32	0.45
3:R:587:LEU:HD22	3:R:644:ALA:HB1	1.99	0.45
2:L:8:PRO:HG3	2:L:11:MET:CE	2.47	0.45
3:R:516:THR:HG22	3:R:600:CYS:SG	2.57	0.45
2:L:139:TYR:CD1	2:L:139:TYR:C	2.86	0.45
2:L:149:ILE:HD12	2:L:154:ARG:HH12	1.81	0.45
1:H:96:CYS:O	1:H:112:GLY:N	2.50	0.45
2:L:93:TYR:O	2:L:93:TYR:CD1	2.70	0.45
3:R:180:LEU:HD21	3:R:182:LEU:HD21	1.98	0.45
3:R:276:ASP:HA	3:R:399:THR:HG1	1.81	0.44
1:H:171:HIS:NE2	2:L:137:ASN:ND2	2.64	0.43
2:L:8:PRO:HG2	2:L:11:MET:HE2	1.98	0.43
2:L:90:TYR:CD1	2:L:95:LEU:HD11	2.53	0.43
3:R:410:ASP:OD1	3:R:418:ARG:NH2	2.52	0.43
1:H:1:GLU:HG2	1:H:2:VAL:HG23	2.01	0.43
1:H:3:GLN:NE2	1:H:25:SER:OG	2.52	0.43
3:R:116:TYR:O	3:R:116:TYR:CG	2.72	0.42
1:H:135:CYS:N	2:L:213:CYS:SG	2.92	0.42



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Atom-1	Atom-2	$egin{aligned} &  ext{Interatomic} \ &  ext{distance} \ &  ext{(Å)} \end{aligned}$	$egin{aligned}  ext{Clash} \  ext{overlap } ( ext{Å}) \end{aligned}$
3:R:111:THR:OG1	3:R:136:THR:HG22	2.20	0.42
3:R:274:ARG:NE	3:R:276:ASP:OD1	2.53	0.41
3:R:219:ASP:OD2	3:R:219:ASP:N	2.53	0.41
2:L:209:ASN:OD1	2:L:210:ARG:N	2.52	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	Perce	entiles
1	Н	$215/217\ (99\%)$	211 (98%)	4 (2%)	0	100	100
2	L	211/213 (99%)	203 (96%)	8 (4%)	0	100	100
3	R	561/702 (80%)	534 (95%)	27 (5%)	0	100	100
All	All	987/1132 (87%)	948 (96%)	39 (4%)	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	Н	187/187 (100%)	186 (100%)	1 (0%)	88 94		
2	L	187/187 (100%)	182 (97%)	5 (3%)	44 74		
3	R	508/636 (80%)	505 (99%)	3 (1%)	86 94		



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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	882/1010 (87%)	873 (99%)	9 (1%)	77 90

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

Mol	Chain	Res	Type
1	Н	77	SER
2	L	7	SER
2	L	93	TYR
2	L	137	ASN
2	L	169	ASP
2	L	171	THR
3	R	35	GLU
3	R	184	LEU
3	R	218	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
2	L	137	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	Trino	Chain	Dag	Link	Bo	ond leng	ths	В	ond ang	les
1	VIOI	Type	Chain	Res	Lilik	Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z >2
	4	NAG	A	1	3,4	14,14,15	0.25	0	17,19,21	0.37	0
	4	NAG	A	2	4	14,14,15	0.20	0	17,19,21	0.45	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
4	NAG	A	1	3,4	-	2/6/23/26	0/1/1/1
4	NAG	A	2	4	-	2/6/23/26	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (4) torsion outliers are listed below:

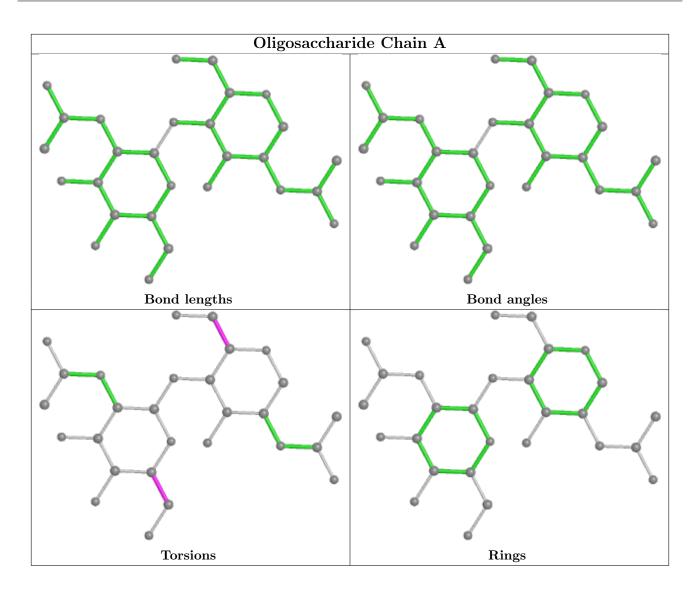
Mol	Chain	Res	Type	Atoms
4	A	2	NAG	O5-C5-C6-O6
4	A	2	NAG	C4-C5-C6-O6
4	A	1	NAG	C4-C5-C6-O6
4	A	1	NAG	O5-C5-C6-O6

There are no ring outliers.

No monomer is involved in short contacts.

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





### 5.6 Ligand geometry (i)

3 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).

Mal	Mol   Type   Ch		Res	Link	Bo	ond leng	$ ag{ths}$	Bond angles		
MIOI	туре	Chain	nes	LIIIK	Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
5	NAG	R	802	3	14,14,15	0.19	0	17,19,21	0.43	0
5	NAG	R	803	3	14,14,15	0.23	0	17,19,21	0.50	0
5	NAG	R	801	3	14,14,15	0.20	0	17,19,21	0.43	0



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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	NAG	R	802	3	-	2/6/23/26	0/1/1/1
5	NAG	R	803	3	-	3/6/23/26	0/1/1/1
5	NAG	R	801	3	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	R	803	NAG	C4-C5-C6-O6
5	R	803	NAG	O5-C5-C6-O6
5	R	802	NAG	O5-C5-C6-O6
5	R	802	NAG	C4-C5-C6-O6
5	R	803	NAG	C3-C2-N2-C7

There are no ring outliers.

No monomer is involved in short contacts.

### 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.



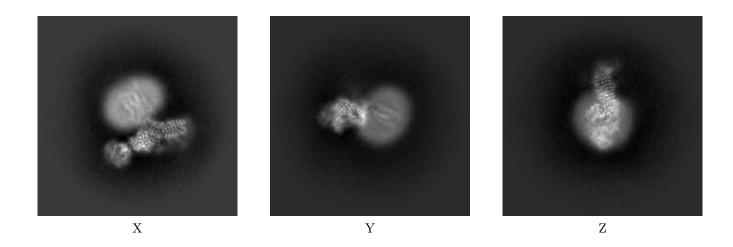
# 6 Map visualisation (i)

This section contains visualisations of the EMDB entry EMD-25762. These allow visual inspection of the internal detail of the map and identification of artifacts.

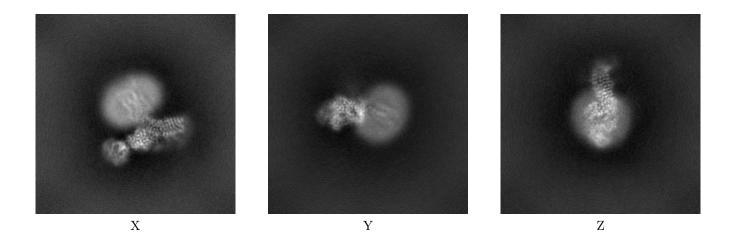
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

#### 6.1 Orthogonal projections (i)

#### 6.1.1 Primary map



#### 6.1.2 Raw map

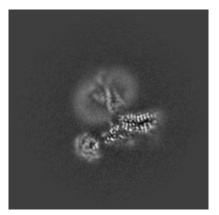


The images above show the map projected in three orthogonal directions.

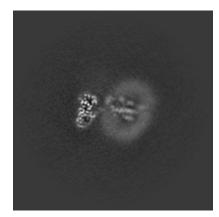


## 6.2 Central slices (i)

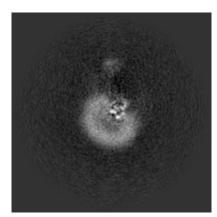
#### 6.2.1 Primary map







Y Index: 224



Z Index: 224

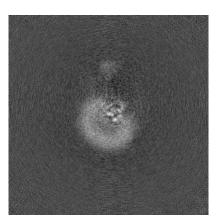
#### 6.2.2 Raw map



X Index: 224



Y Index: 224



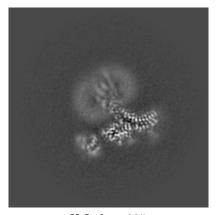
Z Index: 224

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

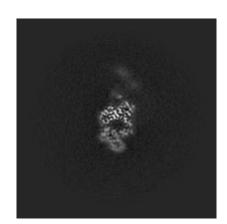


### 6.3 Largest variance slices (i)

#### 6.3.1 Primary map





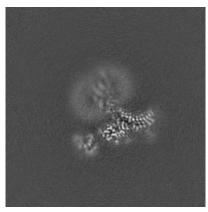


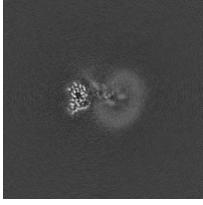
X Index: 235

Y Index: 236

Z Index: 156

#### 6.3.2 Raw map







X Index: 236

Y Index: 236

Z Index: 156

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



The images above show the 3D surface view of the map at the recommended contour level 0.2. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

#### 6.4.2 Raw map



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.



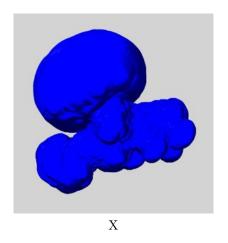
### 6.5 Mask visualisation (i)

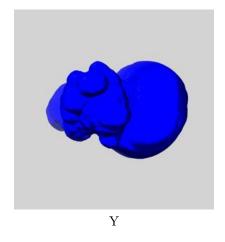
This section shows the 3D surface view of the primary map at 50% transparency overlaid with the specified mask at 0% transparency

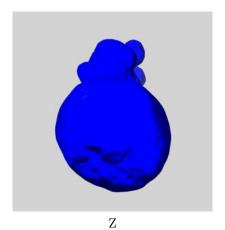
A mask typically either:

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

#### 6.5.1 emd\_25762\_msk\_1.map (i)



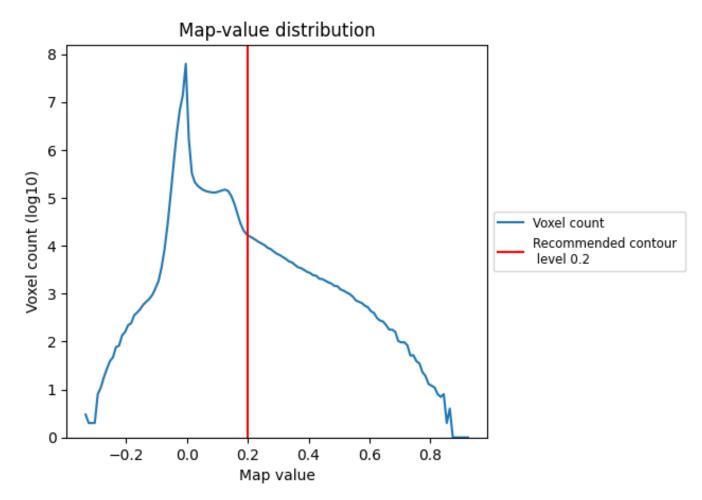




# 7 Map analysis (i)

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

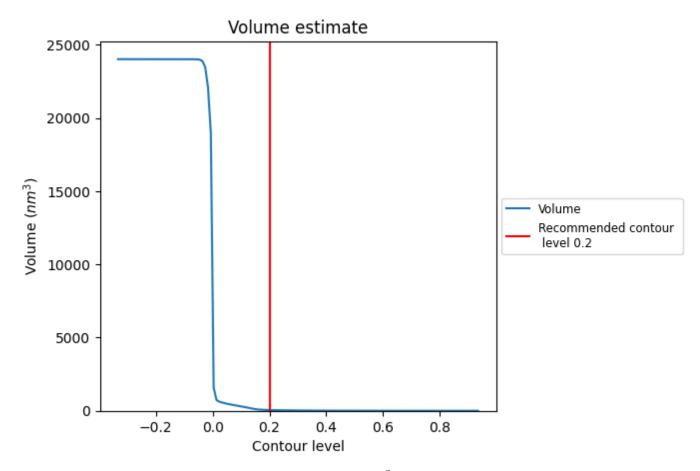
### 7.1 Map-value distribution (i)



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)

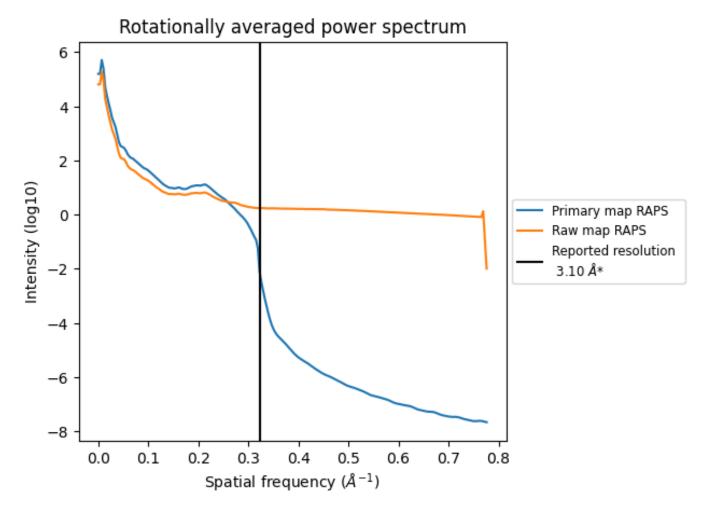


The volume at the recommended contour level is  $52 \text{ nm}^3$ ; this corresponds to an approximate mass of 47 kDa.

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)



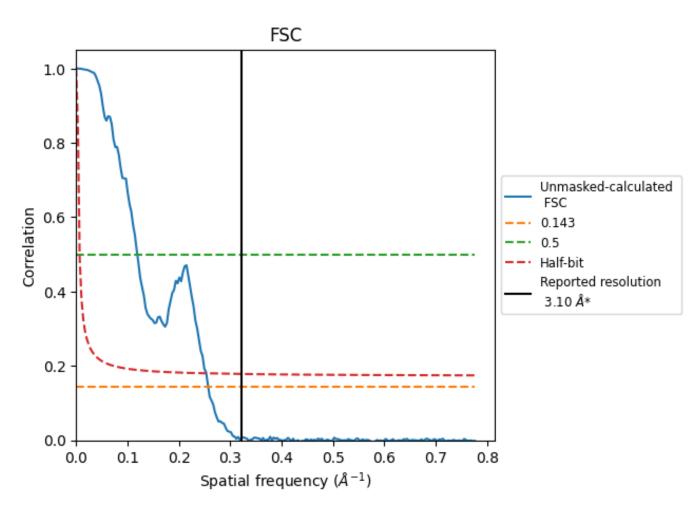
<sup>\*</sup>Reported resolution corresponds to spatial frequency of 0.323  $\rm \mathring{A}^{-1}$ 



# 8 Fourier-Shell correlation (i)

Fourier-Shell Correlation (FSC) is the most commonly used method to estimate the resolution of single-particle and subtomogram-averaged maps. The shape of the curve depends on the imposed symmetry, mask and whether or not the two 3D reconstructions used were processed from a common reference. The reported resolution is shown as a black line. A curve is displayed for the half-bit criterion in addition to lines showing the 0.143 gold standard cut-off and 0.5 cut-off.

#### 8.1 FSC (i)



\*Reported resolution corresponds to spatial frequency of 0.323  $\rm \mathring{A}^{-1}$ 



## 8.2 Resolution estimates (i)

Resolution estimate (Å)	Estimation criterion (FSC cut-off)			
rtesolution estimate (A)	0.143	0.5	Half-bit	
Reported by author	3.10	-	-	
Author-provided FSC curve	-	-	-	
Unmasked-calculated*	3.89	8.40	3.95	

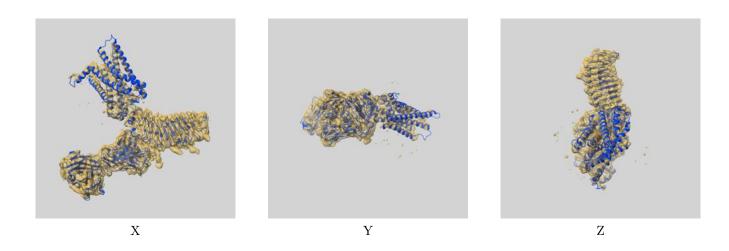
<sup>\*</sup>Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 3.89 differs from the reported value 3.1 by more than 10 %



# 9 Map-model fit (i)

This section contains information regarding the fit between EMDB map EMD-25762 and PDB model 7T9M. Per-residue inclusion information can be found in section 3 on page 6.

### 9.1 Map-model overlay (i)



The images above show the 3D surface view of the map at the recommended contour level 0.2 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

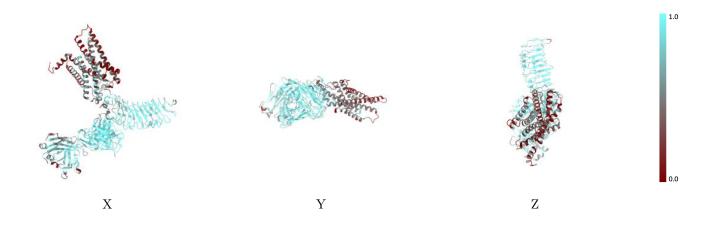


### 9.2 Q-score mapped to coordinate model (i)



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

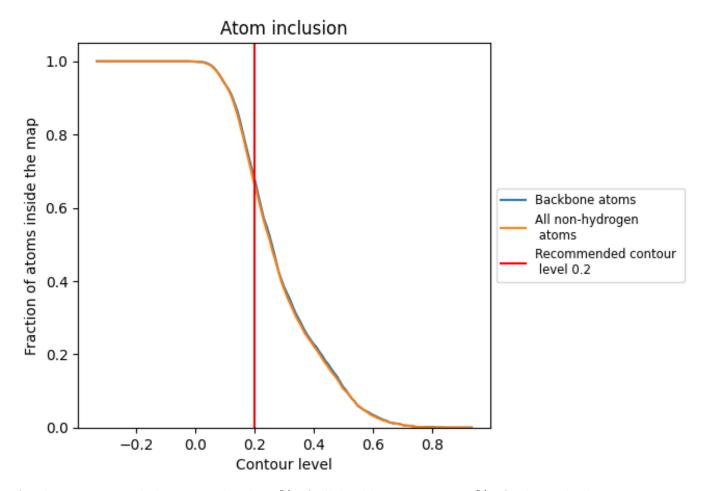
### 9.3 Atom inclusion mapped to coordinate model (i)



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (0.2).



### 9.4 Atom inclusion (i)



At the recommended contour level, 68% 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 (0.2) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.6627	0.3960
A	0.8571	0.5220
Н	0.8104	0.4590
L	0.7931	0.4470
R	0.5693	0.3550



