Full wwPDB EM Validation Report (i)

Oct 30, 2025 – 07:35 PM EDT

PDB ID	:	9YNN / pdb_00009ynn
EMDB ID	:	EMD-73211
Title	:	Human GABAA receptor alpha1-beta2-gamma2 subtype in complex with GABA plus S-nidradine
Deposited on	:	2025-10-10
Resolution	:	2.90 Å(reported)

This wwPDB validation report is for manuscript review

This is a Full wwPDB EM Validation Report.

This report is produced by the wwPDB biocuration pipeline after annotation of the structure.

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:

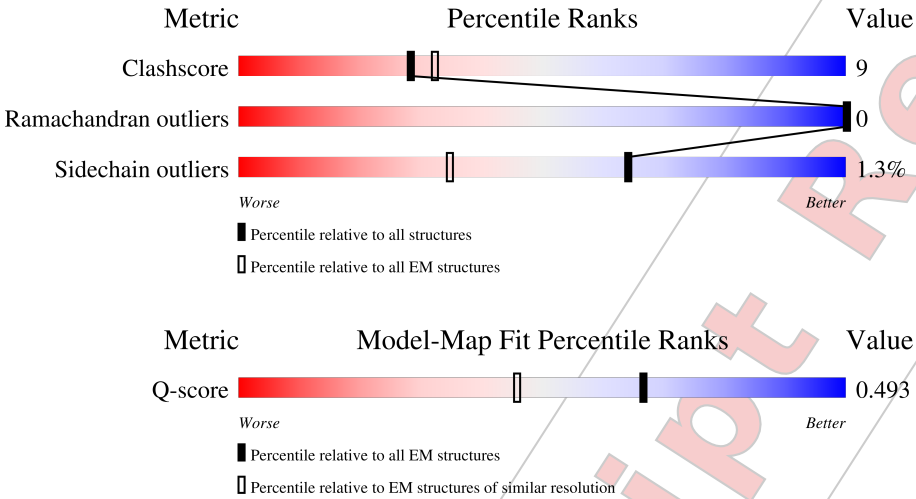
EMDB validation analysis	:	0.0.1.dev129
Mogul	:	2022.3.0, CSD as543be (2022)
MolProbity	:	4-5-2 with Phenix2.0
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
EM percentile statistics	:	202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ	:	1.9.13
Ideal geometry (proteins)	:	Engh & Huber (2001)

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 2.90 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	210492	15764	-
Ramachandran outliers	207382	16835	-
Sidechain outliers	206894	16415	-
Q-score	-	25397	13054 (2.40 - 3.40)

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	B	358	70% 23% • 6%
1	D	358	77% 16% • 6%

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Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.46

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Mol	Chain	Length	Quality of chain
2	A	364	70% 21% 8%
2	C	364	72% 19% 8%
3	E	417	64% 16% 20%
4	I	213	40% 9% 51%
4	L	213	43% 7% 50%
5	J	454	19% 6% 74%
5	K	454	21% 5% 74%
6	F	2	50% 50%
7	G	10	20% 70% 10%
8	H	3	33% 33% 67%
8	M	3	67% 33%

2 Entry composition (i)

There are 12 unique types of molecules in this entry. The entry contains 34796 atoms, of which 17327 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-1.

Mol	Chain	Residues	Atoms						AltConf	Trace
1	B	338	Total	C	H	N	O	S	0	0
			5454	1763	2724	461	490	16		
1	D	338	Total	C	H	N	O	S	0	0
			5454	1763	2724	461	490	16		

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	313	SER	-	linker	UNP P14867
B	314	GLN	-	linker	UNP P14867
B	315	PRO	-	linker	UNP P14867
B	316	ALA	-	linker	UNP P14867
B	317	ARG	-	linker	UNP P14867
B	318	ALA	-	linker	UNP P14867
B	319	ALA	-	linker	UNP P14867
D	313	SER	-	linker	UNP P14867
D	314	GLN	-	linker	UNP P14867
D	315	PRO	-	linker	UNP P14867
D	316	ALA	-	linker	UNP P14867
D	317	ARG	-	linker	UNP P14867
D	318	ALA	-	linker	UNP P14867
D	319	ALA	-	linker	UNP P14867

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

Mol	Chain	Residues	Atoms						AltConf	Trace
2	C	334	Total	C	H	N	O	S	0	0
			5472	1791	2740	440	485	16		
2	A	334	Total	C	H	N	O	S	0	0
			5474	1791	2742	440	485	16		

There are 62 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	308	SER	-	linker	UNP P47870
C	309	GLN	-	linker	UNP P47870
C	310	PRO	-	linker	UNP P47870
C	311	ALA	-	linker	UNP P47870
C	312	ARG	-	linker	UNP P47870
C	313	ALA	-	linker	UNP P47870
C	314	ALA	-	linker	UNP P47870
C	315	ALA	-	linker	UNP P47870
C	342	VAL	-	expression tag	UNP P47870
C	343	ASP	-	expression tag	UNP P47870
C	344	GLY	-	expression tag	UNP P47870
C	345	SER	-	expression tag	UNP P47870
C	346	GLY	-	expression tag	UNP P47870
C	347	ALA	-	expression tag	UNP P47870
C	348	THR	-	expression tag	UNP P47870
C	349	ASN	-	expression tag	UNP P47870
C	350	PHE	-	expression tag	UNP P47870
C	351	SER	-	expression tag	UNP P47870
C	352	LEU	-	expression tag	UNP P47870
C	353	LEU	-	expression tag	UNP P47870
C	354	LYS	-	expression tag	UNP P47870
C	355	GLN	-	expression tag	UNP P47870
C	356	ALA	-	expression tag	UNP P47870
C	357	GLY	-	expression tag	UNP P47870
C	358	ASP	-	expression tag	UNP P47870
C	359	VAL	-	expression tag	UNP P47870
C	360	GLU	-	expression tag	UNP P47870
C	361	GLU	-	expression tag	UNP P47870
C	362	ASN	-	expression tag	UNP P47870
C	363	PRO	-	expression tag	UNP P47870
C	364	GLY	-	expression tag	UNP P47870
A	308	SER	-	linker	UNP P47870
A	309	GLN	-	linker	UNP P47870
A	310	PRO	-	linker	UNP P47870
A	311	ALA	-	linker	UNP P47870
A	312	ARG	-	linker	UNP P47870
A	313	ALA	-	linker	UNP P47870
A	314	ALA	-	linker	UNP P47870
A	315	ALA	-	linker	UNP P47870
A	342	VAL	-	expression tag	UNP P47870
A	343	ASP	-	expression tag	UNP P47870
A	344	GLY	-	expression tag	UNP P47870
A	345	SER	-	expression tag	UNP P47870

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Chain	Residue	Modelled	Actual	Comment	Reference
A	346	GLY	-	expression tag	UNP P47870
A	347	ALA	-	expression tag	UNP P47870
A	348	THR	-	expression tag	UNP P47870
A	349	ASN	-	expression tag	UNP P47870
A	350	PHE	-	expression tag	UNP P47870
A	351	SER	-	expression tag	UNP P47870
A	352	LEU	-	expression tag	UNP P47870
A	353	LEU	-	expression tag	UNP P47870
A	354	LYS	-	expression tag	UNP P47870
A	355	GLN	-	expression tag	UNP P47870
A	356	ALA	-	expression tag	UNP P47870
A	357	GLY	-	expression tag	UNP P47870
A	358	ASP	-	expression tag	UNP P47870
A	359	VAL	-	expression tag	UNP P47870
A	360	GLU	-	expression tag	UNP P47870
A	361	GLU	-	expression tag	UNP P47870
A	362	ASN	-	expression tag	UNP P47870
A	363	PRO	-	expression tag	UNP P47870
A	364	GLY	-	expression tag	UNP P47870

- Molecule 3 is a protein called Isoform 2 of Gamma-aminobutyric acid receptor subunit gam ma-2,Gamma-aminobutyric acid receptor subunit gamma-2.

Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	S		
3	E	333	5442	1781	2713	448	485	15	0	0

There are 66 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	-36	TRP	-	expression tag	UNP P18507
E	-35	SER	-	expression tag	UNP P18507
E	-34	HIS	-	expression tag	UNP P18507
E	-33	PRO	-	expression tag	UNP P18507
E	-32	GLN	-	expression tag	UNP P18507
E	-31	PHE	-	expression tag	UNP P18507
E	-30	GLU	-	expression tag	UNP P18507
E	-29	LYS	-	expression tag	UNP P18507
E	-28	GLY	-	expression tag	UNP P18507
E	-27	GLY	-	expression tag	UNP P18507
E	-26	GLY	-	expression tag	UNP P18507
E	-25	SER	-	expression tag	UNP P18507

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Chain	Residue	Modelled	Actual	Comment	Reference
E	-24	GLY	-	expression tag	UNP P18507
E	-23	GLY	-	expression tag	UNP P18507
E	-22	GLY	-	expression tag	UNP P18507
E	-21	SER	-	expression tag	UNP P18507
E	-20	GLY	-	expression tag	UNP P18507
E	-19	GLY	-	expression tag	UNP P18507
E	-18	SER	-	expression tag	UNP P18507
E	-17	SER	-	expression tag	UNP P18507
E	-16	ALA	-	expression tag	UNP P18507
E	-15	TRP	-	expression tag	UNP P18507
E	-14	SER	-	expression tag	UNP P18507
E	-13	HIS	-	expression tag	UNP P18507
E	-12	PRO	-	expression tag	UNP P18507
E	-11	GLN	-	expression tag	UNP P18507
E	-10	PHE	-	expression tag	UNP P18507
E	-9	GLU	-	expression tag	UNP P18507
E	-8	LYS	-	expression tag	UNP P18507
E	-7	LEU	-	expression tag	UNP P18507
E	-6	GLU	-	expression tag	UNP P18507
E	-5	VAL	-	expression tag	UNP P18507
E	-4	LEU	-	expression tag	UNP P18507
E	-3	PHE	-	expression tag	UNP P18507
E	-2	GLN	-	expression tag	UNP P18507
E	-1	GLY	-	expression tag	UNP P18507
E	0	PRO	-	expression tag	UNP P18507
E	323	SER	-	linker	UNP P18507
E	324	GLN	-	linker	UNP P18507
E	325	PRO	-	linker	UNP P18507
E	326	ALA	-	linker	UNP P18507
E	327	ARG	-	linker	UNP P18507
E	328	ALA	-	linker	UNP P18507
E	358	SER	-	expression tag	UNP P18507
E	359	ARG	-	expression tag	UNP P18507
E	360	GLY	-	expression tag	UNP P18507
E	361	SER	-	expression tag	UNP P18507
E	362	GLY	-	expression tag	UNP P18507
E	363	ALA	-	expression tag	UNP P18507
E	364	THR	-	expression tag	UNP P18507
E	365	ASN	-	expression tag	UNP P18507
E	366	PHE	-	expression tag	UNP P18507
E	367	SER	-	expression tag	UNP P18507
E	368	LEU	-	expression tag	UNP P18507

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Chain	Residue	Modelled	Actual	Comment	Reference
E	369	LEU	-	expression tag	UNP P18507
E	370	LYS	-	expression tag	UNP P18507
E	371	GLN	-	expression tag	UNP P18507
E	372	ALA	-	expression tag	UNP P18507
E	373	GLY	-	expression tag	UNP P18507
E	374	ASP	-	expression tag	UNP P18507
E	375	VAL	-	expression tag	UNP P18507
E	376	GLU	-	expression tag	UNP P18507
E	377	GLU	-	expression tag	UNP P18507
E	378	ASN	-	expression tag	UNP P18507
E	379	PRO	-	expression tag	UNP P18507
E	380	GLY	-	expression tag	UNP P18507

- Molecule 4 is a protein called Kappa FAB light chain.

Mol	Chain	Residues	Atoms						AltConf	Trace
4	I	105	Total	C	H	N	O	S	0	0
			1573	504	771	130	163	5		
4	L	106	Total	C	H	N	O	S	0	0
			1595	510	784	132	164	5		

- Molecule 5 is a protein called IGG2B FAB heavy chain.

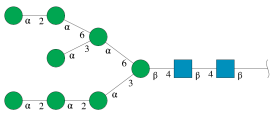
Mol	Chain	Residues	Atoms						AltConf	Trace
5	J	116	Total	C	H	N	O	S	0	0
			1784	574	877	151	178	4		
5	K	117	Total	C	H	N	O	S	0	0
			1798	578	884	152	180	4		

- Molecule 6 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
6	F	2	Total	C	H	N	O	0	0
			53	16	25	2	10		

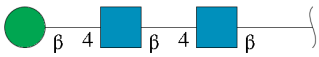
- Molecule 7 is an oligosaccharide called alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-2)-alpha-D-mannopyra

nose-(1-6)-[alpha-D-mannopyranose-(1-3)]alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



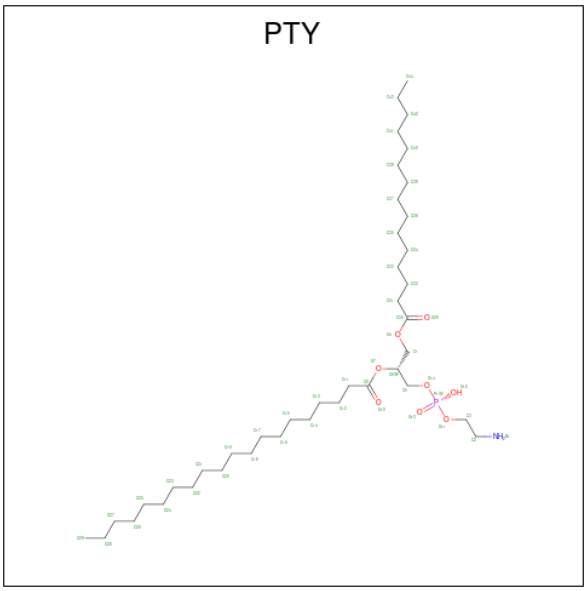
Mol	Chain	Residues	Atoms					AltConf	Trace
7	G	10	Total	C	H	N	O	0	0
			213	64	97	2	50		

- 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.



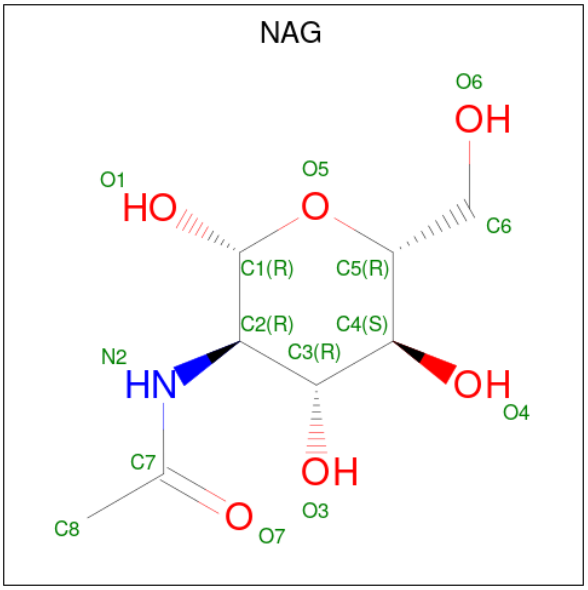
Mol	Chain	Residues	Atoms					AltConf	Trace
8	M	3	Total	C	H	N	O	0	0
			73	22	34	2	15		
8	H	3	Total	C	H	N	O	0	0
			73	22	34	2	15		

- Molecule 9 is PHOSPHATIDYLETHANOLAMINE (CCD ID: PTY) (formula: $C_{40}H_{80}NO_8P$).



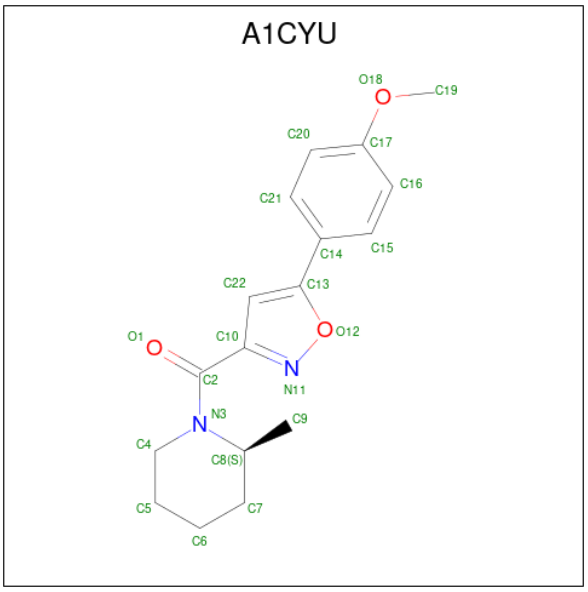
Mol	Chain	Residues	Atoms						AltConf
9	B	1	Total	C	H	N	O	P	0
			99	31	58	1	8	1	
9	C	1	Total	C	H	N	O	P	0
			99	31	58	1	8	1	

- Molecule 10 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).



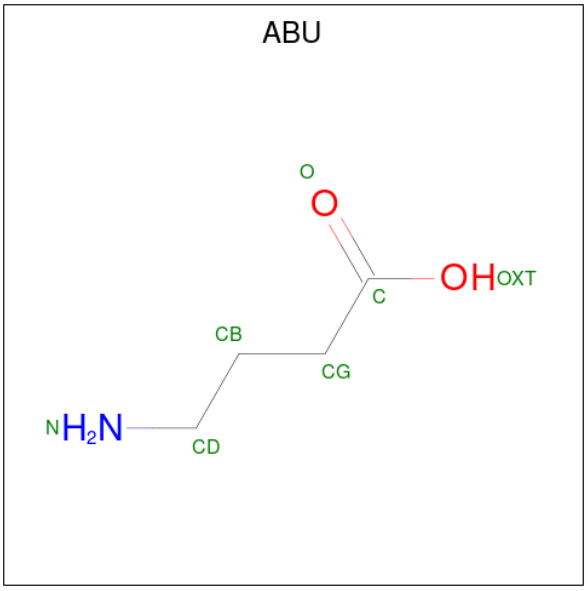
Mol	Chain	Residues	Atoms					AltConf
10	C	1	Total	C	H	N	O	0
			27	8	13	1	5	
10	E	1	Total	C	H	N	O	0
			27	8	13	1	5	
10	A	1	Total	C	N	O		0
			14	8	1	5		

- Molecule 11 is [5-(4-methoxyphenyl)-1,2-oxazol-3-yl][(2S)-2-methylpiperidin-1-yl]methanone (CCD ID: A1CYU) (formula: $C_{17}H_{20}N_2O_3$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					AltConf
11	C	1	Total	C	H	N	O	0
			42	17	20	2	3	

- Molecule 12 is GAMMA-AMINO-BUTANOIC ACID (CCD ID: ABU) (formula: $C_4H_9NO_2$).



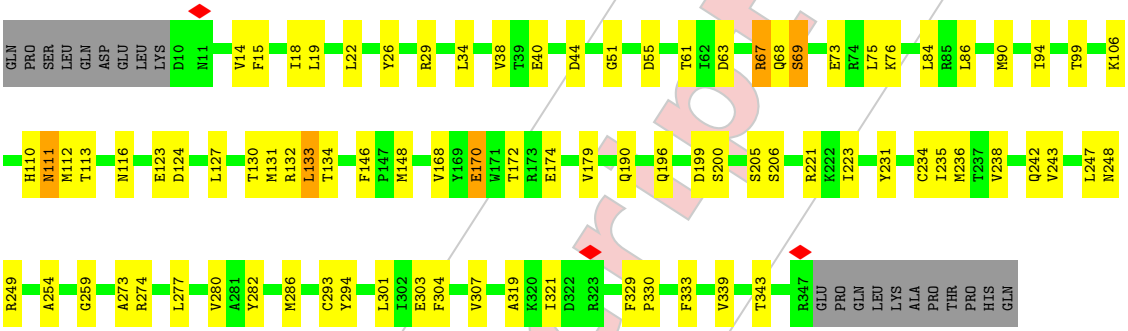
Mol	Chain	Residues	Atoms					AltConf
12	D	1	Total	C	H	N	O	0
			15	4	8	1	2	
12	A	1	Total	C	H	N	O	0
			15	4	8	1	2	

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.

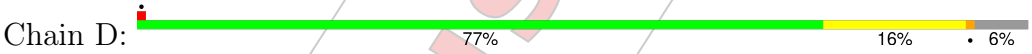
• Molecule 1: Gamma-aminobutyric acid receptor subunit alpha-1

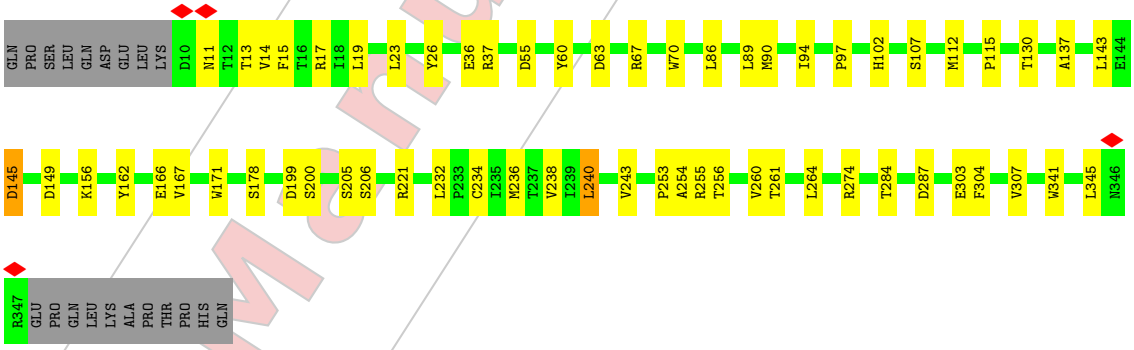
Chain B:



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

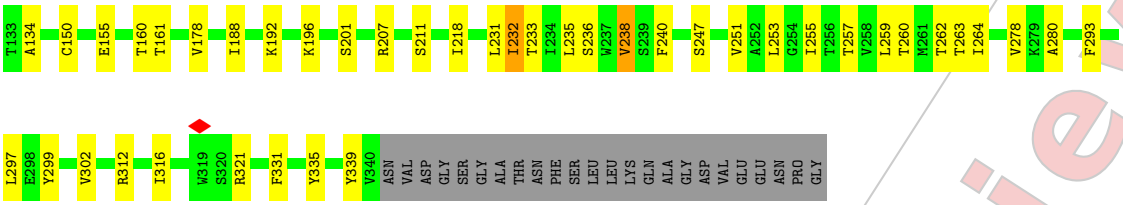
Chain D:





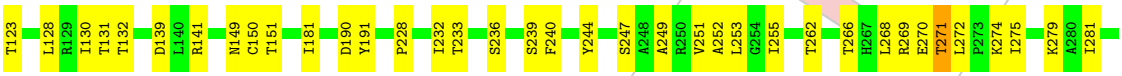
• Molecule 2: Gamma-aminobutyric acid receptor subunit beta-2

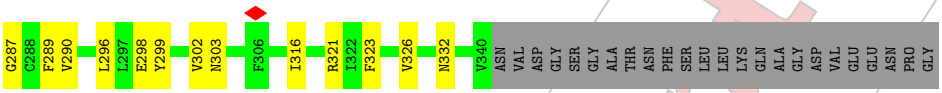
Chain C:



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

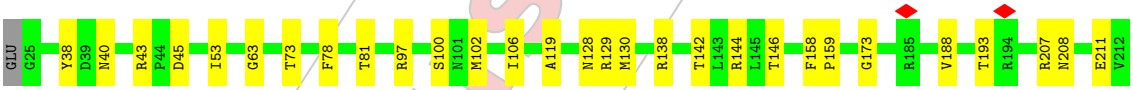
Chain A:

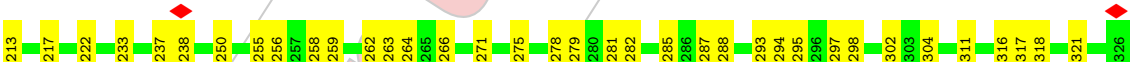


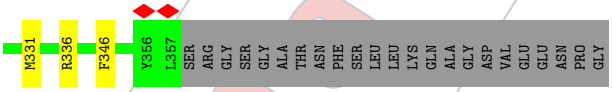


• Molecule 3: Isoform 2 of Gamma-aminobutyric acid receptor subunit gamma-2,Gamma-aminobutyric acid receptor subunit gamma-2

Chain E:

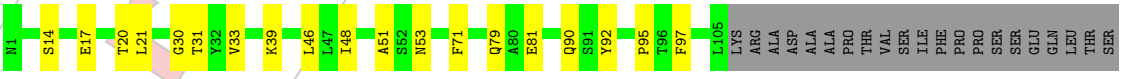


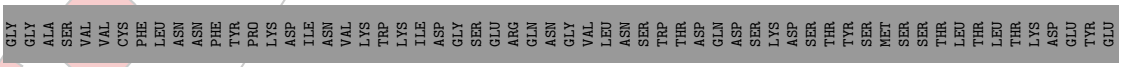


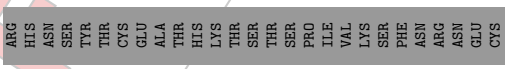


• Molecule 4: Kappa FAB light chain

Chain I:

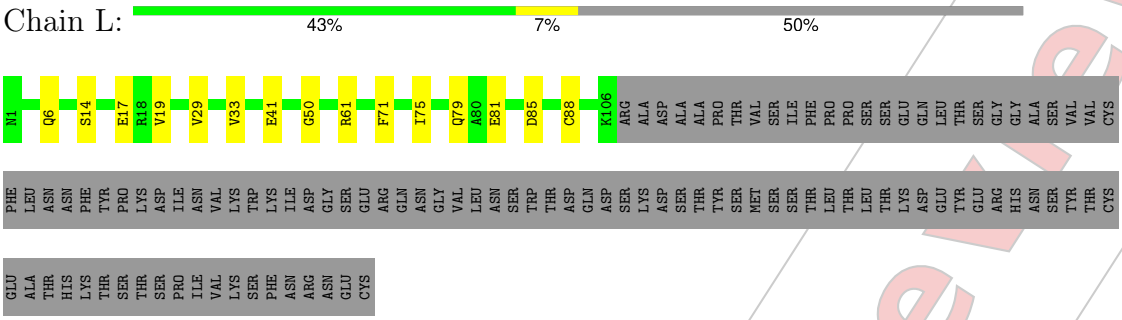






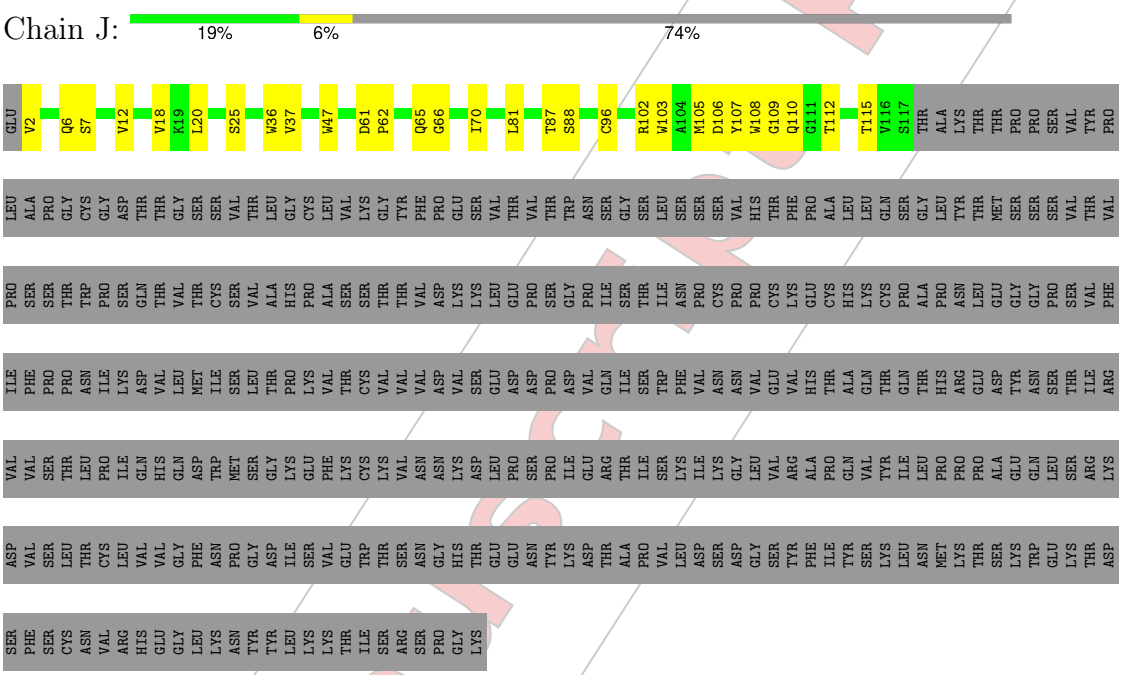
• Molecule 4: Kappa FAB light chain

Chain L:



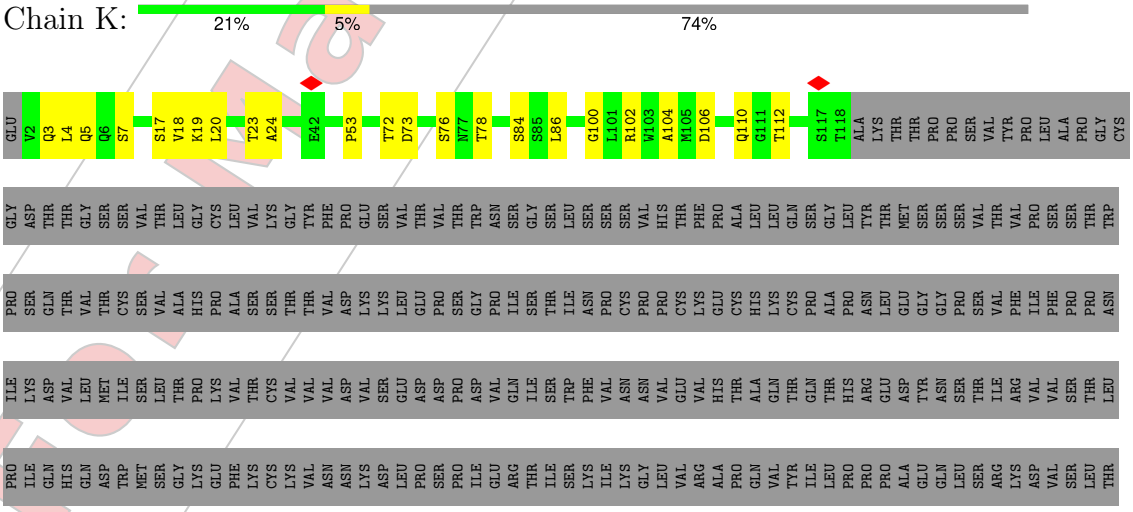
• Molecule 5: IGG2B FAB heavy chain

Chain J:



• Molecule 5: IGG2B FAB heavy chain

Chain K:

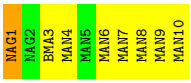


• Molecule 6: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain F:

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

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 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 H:

4 Experimental information (i)

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	64705	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	TFS KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/Å^2$)	50	Depositor
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	2200	Depositor
Magnification	Not provided	
Image detector	FEI FALCON IV (4k x 4k)	Depositor
Maximum map value	2.687	Depositor
Minimum map value	-0.030	Depositor
Average map value	0.002	Depositor
Map value standard deviation	0.031	Depositor
Recommended contour level	0.0297	Depositor
Map size (Å)	269.28, 269.28, 269.28	wwPDB
Map dimensions	288, 288, 288	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.935, 0.935, 0.935	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, PTY, MAN, BMA, A1CYU, ABU

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	B	0.37	2/2799 (0.1%)	0.39	0/3805
1	D	0.32	0/2799	0.38	0/3805
2	A	0.32	0/2804	0.43	0/3818
2	C	0.32	0/2804	0.42	0/3818
3	E	0.28	0/2805	0.39	0/3822
4	I	0.27	0/820	0.34	0/1112
4	L	0.27	0/829	0.36	0/1123
5	J	0.23	0/928	0.34	0/1260
5	K	0.24	0/935	0.33	0/1270
All	All	0.31	2/17523 (0.0%)	0.39	0/23833

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	111	ASN	C-N	6.50	1.42	1.33
1	B	113	THR	C-N	-5.41	1.26	1.33

There are no bond angle outliers.

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	B	2730	2724	2724	63	0
1	D	2730	2724	2724	44	0
2	A	2732	2742	2742	63	0
2	C	2732	2740	2741	55	0
3	E	2729	2713	2714	68	0
4	I	802	771	771	15	0
4	L	811	784	784	11	0
5	J	907	877	877	24	0
5	K	914	884	884	21	0
6	F	28	25	25	1	0
7	G	116	97	97	1	0
8	H	39	34	34	4	0
8	M	39	34	34	1	0
9	B	41	58	58	0	0
9	C	41	58	58	0	0
10	A	14	0	13	1	0
10	C	14	13	13	0	0
10	E	14	13	13	0	0
11	C	22	20	0	0	0
12	A	7	8	0	1	0
12	D	7	8	0	0	0
All	All	17469	17327	17306	324	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 9.

All (324) 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)
5:K:7:SER:O	5:K:112:THR:HG22	1.55	1.06
3:E:317:LEU:O	3:E:321:VAL:HG23	1.66	0.95
3:E:213:VAL:HG22	3:E:222:VAL:HG23	1.51	0.92
5:K:23:THR:OG1	5:K:78:THR:CG2	2.17	0.92
5:K:23:THR:OG1	5:K:78:THR:HG22	1.70	0.90
2:C:113:ASN:OD1	2:C:129:ARG:NH1	2.07	0.88
3:E:81:THR:HG23	3:E:138:ARG:HE	1.39	0.85
3:E:81:THR:CG2	3:E:138:ARG:NE	2.39	0.84
3:E:119:ALA:HB3	2:A:109:VAL:HG11	1.56	0.84
5:J:36:TRP:CD2	5:J:81:LEU:HD22	2.14	0.82
3:E:81:THR:HG23	3:E:138:ARG:NE	1.96	0.81
5:J:36:TRP:CG	5:J:81:LEU:HD22	2.18	0.79
3:E:81:THR:CG2	3:E:138:ARG:HE	1.95	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:L:61:ARG:NH2	4:L:81:GLU:OE2	2.16	0.79
5:J:70:ILE:HG12	5:J:81:LEU:HD13	1.66	0.77
2:C:201:SER:OG	1:D:67:ARG:NH2	2.18	0.77
3:E:302:VAL:HG11	3:E:346:PHE:CZ	2.21	0.76
1:B:67:ARG:NH1	12:A:401:ABU:OXT	2.20	0.75
5:K:23:THR:HA	5:K:78:THR:HG22	1.67	0.74
1:B:190:GLN:NE2	2:A:274:LYS:O	2.22	0.72
1:D:36:GLU:N	1:D:36:GLU:OE1	2.23	0.72
1:D:256:THR:HG21	3:E:264:ALA:HB1	1.71	0.71
3:E:81:THR:HG21	3:E:138:ARG:NE	2.06	0.70
2:C:104:SER:OG	1:D:112:MET:HE3	1.91	0.70
5:K:23:THR:OG1	5:K:78:THR:HG21	1.92	0.69
5:J:2:VAL:N	5:J:25:SER:O	2.27	0.68
5:K:23:THR:CB	5:K:78:THR:HG22	2.23	0.67
1:B:148:MET:HG3	1:B:223:ILE:HD11	1.76	0.67
3:E:213:VAL:CG2	3:E:222:VAL:CG2	2.73	0.67
3:E:213:VAL:CG2	3:E:222:VAL:HG23	2.25	0.67
5:K:23:THR:CA	5:K:78:THR:HG22	2.24	0.66
3:E:213:VAL:HG22	3:E:222:VAL:CG2	2.23	0.66
5:K:3:GLN:NE2	5:K:5:GLN:OE1	2.29	0.65
5:K:7:SER:O	5:K:112:THR:CG2	2.39	0.65
2:A:190:ASP:OD1	2:A:191:TYR:N	2.30	0.65
1:D:274:ARG:NH1	1:D:287:ASP:OD2	2.30	0.64
5:J:96:CYS:O	5:J:109:GLY:N	2.31	0.63
2:C:233:THR:O	2:C:236:SER:OG	2.17	0.62
3:E:233:MET:HA	3:E:233:MET:HE2	1.81	0.62
1:B:123:GLU:N	1:B:123:GLU:OE1	2.28	0.62
5:K:110:GLN:OE1	5:K:110:GLN:N	2.31	0.61
3:E:258:ASN:OD1	3:E:259:LYS:N	2.33	0.61
4:I:79:GLN:OE1	4:I:79:GLN:N	2.33	0.61
1:B:68:GLN:OE1	1:B:99:THR:OG1	2.17	0.61
2:C:196:LYS:NZ	8:H:2:NAG:O7	2.35	0.60
1:D:200:SER:OG	5:K:102:ARG:NH2	2.33	0.60
2:A:84:ASP:OD1	2:A:85:ASN:N	2.35	0.60
1:D:55:ASP:OD2	1:D:221:ARG:NH1	2.33	0.60
3:E:217:SER:O	2:A:117:ARG:NH2	2.35	0.59
3:E:78:PHE:O	3:E:142:THR:HA	2.03	0.59
1:B:40:GLU:OE1	1:B:40:GLU:N	2.36	0.59
3:E:293:VAL:HG13	3:E:297:ASP:OD1	2.04	0.58
1:D:260:VAL:HG11	3:E:271:THR:OG1	2.02	0.58
5:K:4:LEU:HD23	5:K:24:ALA:HB2	1.85	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:L:6:GLN:OE1	4:L:88:CYS:N	2.37	0.57
2:A:302:VAL:HG22	2:A:316:ILE:CD1	2.34	0.57
2:A:323:PHE:O	2:A:326:VAL:HG12	2.04	0.57
1:B:254:ALA:HA	2:A:251:VAL:HG11	1.87	0.57
2:A:271:THR:HG23	2:A:272:LEU:HD12	1.86	0.57
3:E:193:THR:HG22	3:E:193:THR:O	2.04	0.57
5:J:65:GLN:OE1	5:J:66:GLY:N	2.37	0.57
2:A:262:THR:O	2:A:266:THR:HG23	2.05	0.56
2:A:121:ASP:OD2	2:A:123:THR:OG1	2.24	0.56
3:E:173:GLY:O	2:A:117:ARG:NH1	2.39	0.55
4:L:79:GLN:N	4:L:79:GLN:OE1	2.40	0.55
2:A:266:THR:O	2:A:269:ARG:HG2	2.06	0.55
3:E:159:PRO:HG2	3:E:233:MET:HE3	1.89	0.55
5:J:105:MET:O	5:J:108:TRP:NE1	2.38	0.55
2:A:107:HIS:HB2	2:A:113:ASN:OD1	2.06	0.54
3:E:43:ARG:NH1	3:E:45:ASP:O	2.40	0.54
4:L:33:VAL:HG21	4:L:71:PHE:CE1	2.43	0.54
3:E:81:THR:HG23	3:E:138:ARG:CD	2.37	0.54
4:I:14:SER:N	4:I:17:GLU:OE2	2.41	0.54
1:B:179:VAL:O	1:B:196:GLN:NE2	2.40	0.54
1:D:145:ASP:OD1	1:D:145:ASP:N	2.41	0.54
1:D:11:ASN:O	1:D:14:VAL:HG12	2.08	0.53
2:A:79:LEU:HA	10:A:402:NAG:H82	1.91	0.53
2:C:155:GLU:OE1	2:C:207:ARG:NH2	2.41	0.53
2:C:232:ILE:HD12	2:C:260:THR:HG21	1.91	0.53
4:I:97:PHE:CZ	5:J:105:MET:HE1	2.44	0.53
2:A:255:ILE:CD1	2:A:296:LEU:HD21	2.39	0.53
1:B:19:LEU:HD21	1:B:84:LEU:HD13	1.91	0.53
1:B:199:ASP:OD1	1:B:200:SER:N	2.42	0.53
1:B:259:GLY:HA3	1:B:301:LEU:HD13	1.91	0.52
1:D:86:LEU:HD22	1:D:90:MET:SD	2.49	0.52
5:K:73:ASP:OD2	5:K:76:SER:OG	2.25	0.52
2:C:236:SER:HB3	2:C:257:THR:HG21	1.91	0.52
2:A:8:ASN:O	2:A:12:VAL:HG23	2.10	0.52
3:E:316:THR:HG23	3:E:331:MET:HE1	1.92	0.52
2:A:268:LEU:HD21	2:A:281:ILE:HG21	1.92	0.52
2:C:278:VAL:O	2:C:278:VAL:HG13	2.10	0.51
3:E:213:VAL:HG21	3:E:222:VAL:CG2	2.39	0.51
3:E:81:THR:CG2	3:E:138:ARG:CD	2.88	0.51
3:E:63:GLY:N	3:E:73:THR:O	2.43	0.51
2:A:100:ASN:ND2	2:A:151:THR:O	2.41	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:26:TYR:HE2	1:B:75:LEU:HD11	1.74	0.51
2:A:23:TYR:OH	2:A:92:TRP:N	2.41	0.51
2:A:121:ASP:OD1	2:A:121:ASP:N	2.42	0.51
2:A:251:VAL:O	2:A:255:ILE:HG12	2.10	0.51
4:I:39:LYS:NZ	4:I:81:GLU:O	2.42	0.51
1:B:131:MET:HG3	1:B:133:LEU:CD2	2.41	0.50
3:E:294:THR:N	3:E:297:ASP:OD2	2.37	0.50
5:K:100:GLY:N	5:K:104:ALA:O	2.44	0.50
2:A:298:GLU:O	2:A:302:VAL:HG23	2.12	0.50
2:C:293:PHE:CE2	1:D:236:MET:HE3	2.46	0.50
5:K:20:LEU:HD13	5:K:112:THR:OG1	2.10	0.50
1:B:14:VAL:O	1:B:18:ILE:HG12	2.11	0.50
2:A:247:SER:O	2:A:251:VAL:HG12	2.11	0.50
2:C:25:ILE:HG22	2:C:91:LEU:C	2.37	0.50
3:E:213:VAL:HG21	3:E:222:VAL:HG21	1.94	0.50
3:E:256:TRP:HA	3:E:336:ARG:HH21	1.77	0.50
5:J:106:ASP:OD1	5:J:107:TYR:N	2.44	0.50
2:A:149:ASN:OD1	6:F:1:NAG:C1	2.59	0.50
2:A:302:VAL:HG22	2:A:316:ILE:HD11	1.93	0.50
1:B:231:TYR:O	1:B:235:ILE:HG23	2.12	0.49
3:E:318:HIS:HA	3:E:321:VAL:CG2	2.42	0.49
1:D:166:GLU:OE2	3:E:97:ARG:HD2	2.12	0.49
2:C:104:SER:OG	1:D:112:MET:CE	2.59	0.49
2:A:39:GLY:O	2:A:66:ALA:N	2.42	0.49
2:C:79:LEU:HD23	2:C:79:LEU:H	1.78	0.49
3:E:53:ILE:HA	3:E:81:THR:O	2.11	0.49
2:C:259:LEU:HA	2:C:262:THR:HG22	1.94	0.49
1:B:90:MET:HE2	2:A:26:ARG:HD2	1.94	0.49
1:B:248:ASN:H	2:A:303:ASN:HD21	1.61	0.49
1:D:149:ASP:OD1	1:D:149:ASP:N	2.45	0.49
4:I:30:GLY:O	4:I:31:THR:OG1	2.31	0.49
1:B:44:ASP:HB3	1:B:172:THR:HG22	1.95	0.49
1:B:73:GLU:O	1:B:76:LYS:HG3	2.13	0.48
2:A:63:PHE:N	2:A:128:LEU:O	2.38	0.48
1:B:15:PHE:CB	2:A:27:LEU:HD21	2.43	0.48
3:E:237:THR:OG1	3:E:238:ILE:N	2.46	0.48
1:B:55:ASP:OD2	1:B:221:ARG:NH1	2.46	0.48
1:B:199:ASP:OD1	4:I:31:THR:OG1	2.31	0.48
1:D:240:LEU:HA	1:D:243:VAL:HG12	1.95	0.48
4:I:17:GLU:N	4:I:17:GLU:OE1	2.47	0.48
1:B:235:ILE:O	1:B:238:VAL:HG12	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:192:LYS:CE	8:H:1:NAG:O3	2.62	0.48
2:C:211:SER:OG	8:H:1:NAG:O7	2.23	0.48
5:J:37:VAL:HG21	5:J:108:TRP:HZ3	1.79	0.48
2:A:233:THR:O	2:A:236:SER:OG	2.28	0.48
3:E:318:HIS:HA	3:E:321:VAL:HG23	1.94	0.47
4:L:19:VAL:HG12	4:L:75:ILE:HB	1.95	0.47
3:E:213:VAL:CG2	3:E:222:VAL:HG21	2.44	0.47
1:B:112:MET:HE2	2:A:130:ILE:HG23	1.97	0.47
2:C:115:MET:HE3	2:C:117:ARG:HB2	1.95	0.47
1:B:67:ARG:HG2	1:B:130:THR:HG23	1.95	0.47
3:E:262:VAL:HG12	2:A:249:ALA:CB	2.45	0.47
2:A:287:GLY:O	2:A:290:VAL:HG12	2.15	0.47
2:C:58:THR:HA	2:C:132:THR:O	2.15	0.47
2:C:260:THR:O	2:C:263:THR:HG22	2.15	0.47
3:E:130:MET:HE2	3:E:142:THR:HB	1.97	0.47
4:I:31:THR:HA	4:I:51:ALA:HB2	1.97	0.47
2:A:247:SER:HA	2:A:299:TYR:OH	2.14	0.47
5:J:20:LEU:HD12	5:J:112:THR:HG21	1.96	0.46
1:B:15:PHE:HB2	2:A:27:LEU:HD21	1.96	0.46
2:C:23:TYR:CD1	2:C:25:ILE:HG23	2.49	0.46
2:A:58:THR:HA	2:A:132:THR:O	2.14	0.46
2:A:240:PHE:HB3	2:A:321:ARG:HG2	1.97	0.46
1:D:102:HIS:N	1:D:156:LYS:O	2.49	0.46
3:E:81:THR:HG23	3:E:138:ARG:HD2	1.97	0.46
4:I:95:PRO:HG2	5:J:47:TRP:CD1	2.51	0.46
2:A:228:PRO:O	2:A:232:ILE:HG12	2.16	0.46
1:B:243:VAL:O	1:B:247:LEU:HD23	2.16	0.46
1:B:274:ARG:NH2	1:B:280:VAL:HB	2.31	0.46
1:B:339:VAL:O	1:B:343:THR:HG23	2.16	0.46
1:D:255:ARG:HD2	1:D:304:PHE:CE1	2.51	0.46
4:I:33:VAL:HG21	4:I:71:PHE:CD2	2.51	0.46
2:C:11:LEU:HD11	2:C:76:VAL:HG11	1.97	0.46
2:A:44:ILE:CG2	2:A:181:ILE:HD11	2.45	0.46
2:C:25:ILE:HD12	2:C:26:ARG:N	2.30	0.45
2:C:247:SER:HA	2:C:299:TYR:OH	2.16	0.45
1:B:146:PHE:CD2	1:B:277:LEU:HD11	2.51	0.45
2:C:53:VAL:HG23	2:C:54:ASN:N	2.32	0.45
2:C:121:ASP:OD1	2:C:122:GLY:N	2.49	0.45
2:C:160:THR:HG22	2:C:161:THR:N	2.31	0.45
1:D:36:GLU:O	1:D:37:ARG:HB2	2.15	0.45
5:J:7:SER:O	5:J:112:THR:HG22	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:273:ALA:HB1	1:B:286:MET:HE3	1.99	0.45
1:D:11:ASN:O	1:D:15:PHE:HD1	1.98	0.45
4:I:48:ILE:HG23	4:I:53:ASN:H	1.81	0.45
1:B:26:TYR:CE1	1:B:94:ILE:HA	2.51	0.45
1:D:234:CYS:O	1:D:238:VAL:HG23	2.17	0.45
3:E:173:GLY:O	2:A:82:THR:HG21	2.17	0.45
3:E:263:PRO:HA	3:E:266:THR:OG1	2.16	0.45
1:B:294:TYR:CE1	2:C:231:LEU:HD13	2.51	0.45
1:D:199:ASP:OD1	1:D:200:SER:N	2.49	0.45
1:B:38:VAL:HG11	1:B:168:VAL:HG23	1.98	0.45
2:C:57:TYR:CE2	2:C:150:CYS:HB3	2.52	0.45
3:E:211:GLU:HG2	3:E:222:VAL:HB	1.98	0.45
1:B:236:MET:CE	2:A:289:PHE:HE1	2.29	0.45
2:A:269:ARG:HG3	2:A:270:GLU:N	2.30	0.45
1:D:60:TYR:CE1	1:D:137:ALA:HB3	2.52	0.45
3:E:159:PRO:HB2	3:E:233:MET:HE3	1.99	0.45
2:C:63:PHE:O	2:C:127:GLY:HA2	2.17	0.45
2:C:259:LEU:O	2:C:262:THR:HG22	2.17	0.45
5:J:61:ASP:OD1	5:J:62:PRO:HD2	2.17	0.45
1:B:124:ASP:OD1	1:B:124:ASP:N	2.50	0.44
1:B:280:VAL:HG12	1:B:282:TYR:H	1.80	0.44
1:B:303:GLU:OE1	1:B:304:PHE:N	2.50	0.44
5:K:86:LEU:HD23	5:K:86:LEU:H	1.83	0.44
5:K:106:ASP:OD1	5:K:106:ASP:N	2.50	0.44
3:E:287:LEU:HD12	3:E:288:PRO:O	2.18	0.44
3:E:298:LEU:O	3:E:302:VAL:HG12	2.17	0.44
5:J:12:VAL:HG11	5:J:18:VAL:HG12	1.99	0.44
5:J:36:TRP:CD2	5:J:81:LEU:CD2	2.93	0.44
3:E:193:THR:O	3:E:193:THR:CG2	2.66	0.44
3:E:217:SER:HB2	2:A:117:ARG:HH22	1.82	0.44
4:I:20:THR:C	4:I:21:LEU:HD12	2.42	0.44
4:L:29:VAL:HG23	4:L:29:VAL:O	2.18	0.44
1:D:143:LEU:O	1:D:284:THR:HG22	2.18	0.44
5:J:6:GLN:N	5:J:110:GLN:OE1	2.43	0.44
1:B:170:GLU:OE1	5:J:103:TRP:NE1	2.49	0.44
1:D:15:PHE:O	1:D:19:LEU:HD13	2.18	0.44
5:J:36:TRP:CE3	5:J:81:LEU:CD2	3.01	0.44
2:A:27:LEU:HD23	2:A:31:PHE:CG	2.52	0.44
3:E:217:SER:HB2	2:A:117:ARG:NH2	2.32	0.44
2:A:57:TYR:CE1	2:A:150:CYS:HB3	2.53	0.44
2:C:218:ILE:HD12	2:C:218:ILE:C	2.42	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:51:GLY:N	1:B:61:THR:O	2.49	0.44
1:B:106:LYS:NZ	2:C:105:PHE:CZ	2.84	0.44
1:B:238:VAL:HG11	1:B:333:PHE:CZ	2.53	0.44
2:C:235:LEU:O	2:C:238:VAL:HG22	2.18	0.44
2:C:293:PHE:CE2	1:D:240:LEU:CD2	3.01	0.44
3:E:40:ASN:OD1	3:E:40:ASN:N	2.51	0.44
3:E:100:SER:HB3	3:E:129:ARG:HB2	2.00	0.44
2:A:289:PHE:CD1	2:A:289:PHE:C	2.95	0.44
1:B:274:ARG:O	1:B:274:ARG:CG	2.66	0.43
2:C:280:ALA:HB2	2:C:339:TYR:CE1	2.53	0.43
1:B:134:THR:O	1:B:134:THR:HG23	2.18	0.43
1:D:162:TYR:HB3	1:D:167:VAL:HG23	2.01	0.43
1:D:232:LEU:O	1:D:236:MET:HG2	2.19	0.43
3:E:287:LEU:HB2	3:E:288:PRO:HD2	1.99	0.43
1:B:205:SER:OG	1:B:206:SER:N	2.51	0.43
1:D:205:SER:OG	1:D:206:SER:N	2.51	0.43
3:E:146:THR:O	3:E:146:THR:HG23	2.18	0.43
3:E:304:PHE:HD1	3:E:304:PHE:O	2.02	0.43
2:A:20:LEU:HD11	2:A:91:LEU:HD21	2.00	0.43
2:C:178:VAL:HG12	2:C:188:ILE:HG21	2.00	0.43
1:D:13:THR:O	1:D:17:ARG:HG3	2.19	0.43
1:B:29:ARG:HG2	2:C:87:VAL:HG23	2.01	0.43
4:L:61:ARG:NE	4:L:79:GLN:OE1	2.52	0.43
1:B:238:VAL:O	1:B:242:GLN:HG2	2.18	0.43
2:A:275:ILE:HD11	2:A:279:LYS:HD3	2.00	0.43
1:B:174:GLU:OE1	1:B:174:GLU:N	2.47	0.43
2:C:255:ILE:HG21	1:D:261:THR:HG21	2.01	0.43
4:I:31:THR:O	5:J:102:ARG:NH1	2.52	0.43
5:J:87:THR:HG22	5:J:88:SER:H	1.83	0.43
2:C:101:ASP:HA	2:C:134:ALA:HA	2.01	0.43
2:C:116:ILE:HG23	2:C:116:ILE:O	2.19	0.43
2:C:293:PHE:O	2:C:297:LEU:HG	2.19	0.42
2:A:36:VAL:HG23	2:A:36:VAL:O	2.18	0.42
1:B:15:PHE:HA	1:B:18:ILE:HG12	2.00	0.42
2:C:25:ILE:HD13	1:D:89:LEU:HD13	2.00	0.42
2:C:299:TYR:HA	2:C:302:VAL:HG22	2.02	0.42
1:D:23:LEU:HD21	1:D:94:ILE:HD13	2.00	0.42
5:J:37:VAL:HG21	5:J:108:TRP:CZ3	2.55	0.42
2:A:15:THR:O	2:A:19:LEU:HG	2.19	0.42
5:J:36:TRP:CE3	5:J:81:LEU:HD22	2.52	0.42
4:L:41:GLU:CG	4:L:41:GLU:O	2.68	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:132:ARG:NH2	2:A:101:ASP:OD1	2.50	0.42
2:C:257:THR:HA	2:C:260:THR:HG22	2.01	0.42
1:D:171:TRP:CE3	1:D:178:SER:HB3	2.55	0.42
4:L:85:ASP:OD1	4:L:85:ASP:N	2.53	0.42
3:E:158:PHE:O	3:E:295:ALA:HB3	2.20	0.42
3:E:278:THR:O	3:E:281:THR:OG1	2.32	0.42
5:K:17:SER:OG	5:K:84:SER:O	2.35	0.42
1:B:69:SER:OG	1:B:127:LEU:O	2.30	0.42
1:B:133:LEU:HD13	2:C:109:VAL:HG11	2.02	0.42
2:C:299:TYR:O	2:C:302:VAL:HG22	2.20	0.42
2:A:63:PHE:HB2	2:A:130:ILE:HD12	2.01	0.42
2:A:266:THR:HA	2:A:269:ARG:HG2	2.01	0.42
3:E:250:LEU:C	3:E:250:LEU:HD13	2.45	0.42
3:E:302:VAL:HG11	3:E:346:PHE:CE2	2.55	0.42
5:J:36:TRP:CE2	5:J:81:LEU:HB2	2.55	0.42
4:L:14:SER:N	4:L:17:GLU:OE2	2.53	0.42
1:B:111:ASN:O	7:G:1:NAG:H82	2.20	0.41
1:D:26:TYR:CD1	1:D:26:TYR:C	2.98	0.41
2:A:118:LEU:HA	2:A:123:THR:O	2.20	0.41
1:B:86:LEU:HD13	1:B:90:MET:HG3	2.01	0.41
1:B:236:MET:HE3	2:A:289:PHE:HE1	1.86	0.41
1:B:307:VAL:HG22	1:B:321:ILE:CG2	2.50	0.41
4:L:33:VAL:O	4:L:50:GLY:O	2.39	0.41
1:D:115:PRO:HD3	8:M:1:NAG:H62	2.02	0.41
3:E:128:ASN:ND2	3:E:144:ARG:HH11	2.18	0.41
3:E:255:PHE:O	3:E:336:ARG:NH2	2.53	0.41
3:E:262:VAL:N	3:E:263:PRO:HD2	2.36	0.41
3:E:311:LEU:HD13	3:E:311:LEU:C	2.46	0.41
5:K:4:LEU:CD2	5:K:24:ALA:HB2	2.49	0.41
1:B:75:LEU:HD12	1:B:75:LEU:N	2.36	0.41
3:E:159:PRO:CG	3:E:233:MET:HE3	2.50	0.41
2:A:139:ASP:OD1	2:A:141:ARG:HG3	2.21	0.41
1:B:110:HIS:HB2	1:B:116:ASN:ND2	2.35	0.41
1:B:234:CYS:HB3	1:B:293:CYS:SG	2.60	0.41
2:C:192:LYS:HE2	8:H:1:NAG:O3	2.21	0.41
3:E:38:TYR:CE1	3:E:106:ILE:HA	2.56	0.41
5:K:53:PRO:HA	5:K:72:THR:HG21	2.03	0.41
1:B:249:ARG:NH2	1:B:319:ALA:CB	2.84	0.41
2:C:253:LEU:O	2:C:257:THR:OG1	2.34	0.41
2:A:251:VAL:HG13	2:A:252:ALA:N	2.35	0.41
2:C:251:VAL:HG21	1:D:254:ALA:HB1	2.02	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:253:PRO:HA	1:D:256:THR:OG1	2.21	0.40
1:D:341:TRP:O	1:D:345:LEU:HB2	2.21	0.40
4:I:46:LEU:C	4:I:46:LEU:HD13	2.46	0.40
2:A:100:ASN:OD1	2:A:100:ASN:N	2.50	0.40
2:C:240:PHE:HB3	2:C:321:ARG:HG2	2.04	0.40
1:D:70:TRP:CH2	1:D:97:PRO:HD3	2.56	0.40
1:D:303:GLU:O	1:D:307:VAL:HG23	2.21	0.40
2:C:331:PHE:CD1	2:C:331:PHE:C	2.98	0.40
3:E:188:VAL:HG21	3:E:208:ASN:OD1	2.21	0.40
4:I:90:GLN:OE1	4:I:92:TYR:N	2.54	0.40
2:A:130:ILE:HG22	2:A:131:THR:N	2.36	0.40
2:A:239:SER:HB3	2:A:253:LEU:HD22	2.04	0.40
1:B:116:ASN:ND2	1:B:132:ARG:HE	2.19	0.40
2:C:312:ARG:O	2:C:316:ILE:HG12	2.21	0.40
2:C:335:TYR:CD1	2:C:335:TYR:C	3.00	0.40
1:D:264:LEU:HD11	3:E:275:THR:OG1	2.21	0.40
3:E:102:MET:O	3:E:102:MET:CG	2.69	0.40
1:B:34:LEU:HB2	2:C:9:MET:HE1	2.03	0.40
1:B:329:PHE:HB2	1:B:330:PRO:HD3	2.03	0.40
2:C:59:LEU:O	2:C:131:THR:HA	2.21	0.40
1:D:70:TRP:CH2	1:D:97:PRO:CD	3.05	0.40
1:D:89:LEU:HD23	1:D:89:LEU:O	2.21	0.40
3:E:279:LEU:HA	3:E:282:ILE:HG22	2.02	0.40
5:K:18:VAL:CG1	5:K:19:LYS:N	2.85	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	B	336/358 (94%)	328 (98%)	8 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	D	336/358 (94%)	324 (96%)	12 (4%)	0	100	100
2	A	332/364 (91%)	321 (97%)	11 (3%)	0	100	100
2	C	332/364 (91%)	317 (96%)	15 (4%)	0	100	100
3	E	331/417 (79%)	312 (94%)	19 (6%)	0	100	100
4	I	103/213 (48%)	93 (90%)	10 (10%)	0	100	100
4	L	104/213 (49%)	96 (92%)	8 (8%)	0	100	100
5	J	114/454 (25%)	109 (96%)	5 (4%)	0	100	100
5	K	115/454 (25%)	104 (90%)	11 (10%)	0	100	100
All	All	2103/3195 (66%)	2004 (95%)	99 (5%)	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	B	300/319 (94%)	294 (98%)	6 (2%)	50	79
1	D	300/319 (94%)	295 (98%)	5 (2%)	56	83
2	A	302/326 (93%)	298 (99%)	4 (1%)	65	88
2	C	302/326 (93%)	296 (98%)	6 (2%)	50	79
3	E	305/372 (82%)	303 (99%)	2 (1%)	81	94
4	I	89/188 (47%)	89 (100%)	0	100	100
4	L	90/188 (48%)	90 (100%)	0	100	100
5	J	97/407 (24%)	96 (99%)	1 (1%)	73	91
5	K	98/407 (24%)	98 (100%)	0	100	100
All	All	1883/2852 (66%)	1859 (99%)	24 (1%)	64	88

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	22	LEU
1	B	63	ASP
1	B	67	ARG
1	B	69	SER
1	B	133	LEU
1	B	170	GLU
2	C	25	ILE
2	C	40	MET
2	C	125	LEU
2	C	232	ILE
2	C	238	VAL
2	C	264	ILE
1	D	63	ASP
1	D	107	SER
1	D	130	THR
1	D	145	ASP
1	D	240	LEU
3	E	207	ARG
3	E	285	LYS
5	J	115	THR
2	A	79	LEU
2	A	244	TYR
2	A	271	THR
2	A	332	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	116	ASN
1	B	216	HIS
1	B	248	ASN
1	B	346	ASN
2	C	54	ASN
2	C	148	GLN
2	C	185	GLN
2	C	267	HIS
1	D	56	HIS
1	D	314	GLN
3	E	33	ASN
3	E	60	ASN
3	E	128	ASN
3	E	154	GLN
4	L	1	ASN

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Mol	Chain	Res	Type
2	A	54	ASN
2	A	64	GLN
2	A	303	ASN

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

18 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	$\#\|Z\| > 2$
6	NAG	F	1	6	14,14,15	0.93	1 (7%)	17,19,21	0.55	0
6	NAG	F	2	6	14,14,15	0.56	0	17,19,21	0.45	0
7	NAG	G	1	7,1	14,14,15	0.62	1 (7%)	17,19,21	0.48	0
7	MAN	G	10	7	11,11,12	0.69	0	15,15,17	1.17	2 (13%)
7	NAG	G	2	7	14,14,15	0.19	0	17,19,21	0.52	0
7	BMA	G	3	7	11,11,12	0.79	0	15,15,17	1.06	1 (6%)
7	MAN	G	4	7	11,11,12	0.51	0	15,15,17	1.43	1 (6%)
7	MAN	G	5	7	11,11,12	0.82	0	15,15,17	0.85	0
7	MAN	G	6	7	11,11,12	0.60	0	15,15,17	0.93	2 (13%)
7	MAN	G	7	7	11,11,12	0.66	0	15,15,17	1.22	2 (13%)
7	MAN	G	8	7	11,11,12	0.79	1 (9%)	15,15,17	0.86	1 (6%)
7	MAN	G	9	7	11,11,12	0.75	0	15,15,17	0.91	1 (6%)
8	NAG	H	1	8,2	14,14,15	0.65	1 (7%)	17,19,21	0.45	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
8	NAG	H	2	8	14,14,15	0.63	1 (7%)	17,19,21	0.68	1 (5%)
8	BMA	H	3	8	11,11,12	0.64	0	15,15,17	0.75	0
8	NAG	M	1	8,1	14,14,15	0.48	0	17,19,21	0.43	0
8	NAG	M	2	8	14,14,15	0.21	0	17,19,21	0.63	0
8	BMA	M	3	8	11,11,12	0.60	0	15,15,17	0.82	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
6	NAG	F	1	6	-	0/6/23/26	0/1/1/1
6	NAG	F	2	6	-	4/6/23/26	0/1/1/1
7	NAG	G	1	7,1	-	2/6/23/26	0/1/1/1
7	MAN	G	10	7	-	1/2/19/22	0/1/1/1
7	NAG	G	2	7	-	1/6/23/26	0/1/1/1
7	BMA	G	3	7	-	0/2/19/22	0/1/1/1
7	MAN	G	4	7	-	0/2/19/22	0/1/1/1
7	MAN	G	5	7	-	1/2/19/22	0/1/1/1
7	MAN	G	6	7	-	0/2/19/22	0/1/1/1
7	MAN	G	7	7	-	0/2/19/22	0/1/1/1
7	MAN	G	8	7	-	0/2/19/22	0/1/1/1
7	MAN	G	9	7	-	0/2/19/22	0/1/1/1
8	NAG	H	1	8,2	-	2/6/23/26	0/1/1/1
8	NAG	H	2	8	-	0/6/23/26	0/1/1/1
8	BMA	H	3	8	-	1/2/19/22	0/1/1/1
8	NAG	M	1	8,1	-	1/6/23/26	0/1/1/1
8	NAG	M	2	8	-	0/6/23/26	0/1/1/1
8	BMA	M	3	8	-	0/2/19/22	0/1/1/1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	F	1	NAG	O5-C1	-3.34	1.38	1.43
8	H	1	NAG	O5-C1	-2.25	1.39	1.43
8	H	2	NAG	O5-C1	-2.11	1.40	1.43
7	G	1	NAG	O5-C1	-2.09	1.40	1.43
7	G	8	MAN	O5-C1	-2.03	1.40	1.43

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
7	G	4	MAN	C1-O5-C5	4.47	118.17	112.19
7	G	10	MAN	C1-O5-C5	3.11	116.35	112.19
7	G	7	MAN	O2-C2-C3	-2.90	104.14	110.15
8	H	2	NAG	C1-O5-C5	2.51	115.55	112.19
7	G	7	MAN	C1-O5-C5	2.40	115.40	112.19
7	G	10	MAN	O2-C2-C3	-2.28	105.43	110.15
7	G	9	MAN	O2-C2-C3	-2.20	105.59	110.15
7	G	8	MAN	O2-C2-C3	-2.14	105.73	110.15
7	G	6	MAN	O2-C2-C3	-2.10	105.80	110.15
7	G	3	BMA	O2-C2-C3	-2.03	105.94	110.15
7	G	6	MAN	C1-O5-C5	2.01	114.88	112.19

There are no chirality outliers.

All (13) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	F	2	NAG	C1-C2-N2-C7
7	G	1	NAG	O5-C5-C6-O6
7	G	1	NAG	C4-C5-C6-O6
7	G	5	MAN	O5-C5-C6-O6
8	H	3	BMA	O5-C5-C6-O6
7	G	10	MAN	O5-C5-C6-O6
6	F	2	NAG	C4-C5-C6-O6
8	H	1	NAG	C3-C2-N2-C7
6	F	2	NAG	O5-C5-C6-O6
7	G	2	NAG	C1-C2-N2-C7
8	M	1	NAG	C1-C2-N2-C7
8	H	1	NAG	C1-C2-N2-C7
6	F	2	NAG	C3-C2-N2-C7

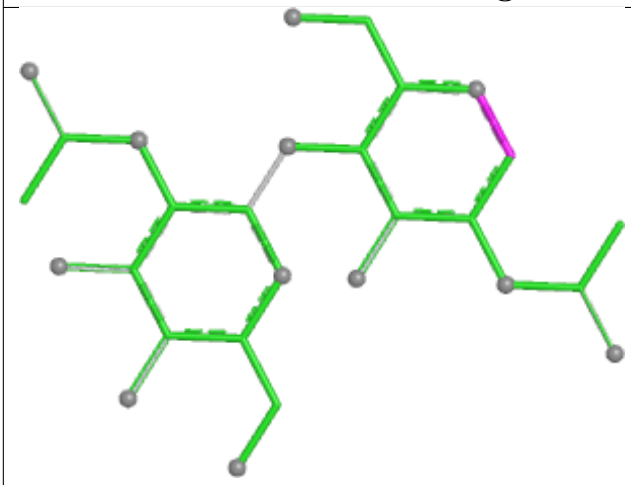
There are no ring outliers.

5 monomers are involved in 7 short contacts:

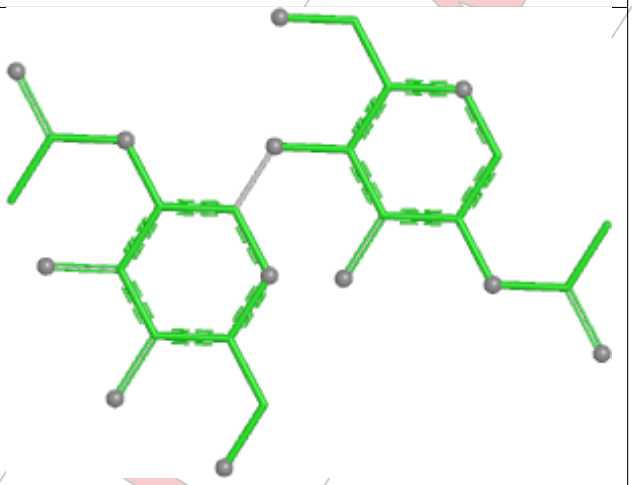
Mol	Chain	Res	Type	Clashes	Symm-Clashes
8	H	1	NAG	3	0
8	M	1	NAG	1	0
8	H	2	NAG	1	0
7	G	1	NAG	1	0
6	F	1	NAG	1	0

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

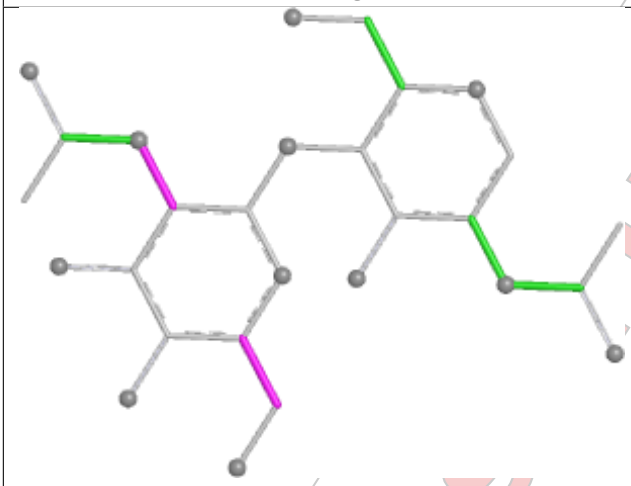
Oligosaccharide Chain F



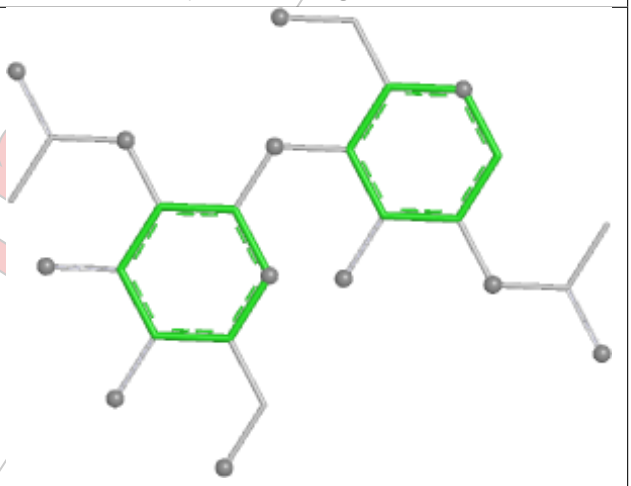
Bond lengths



Bond angles

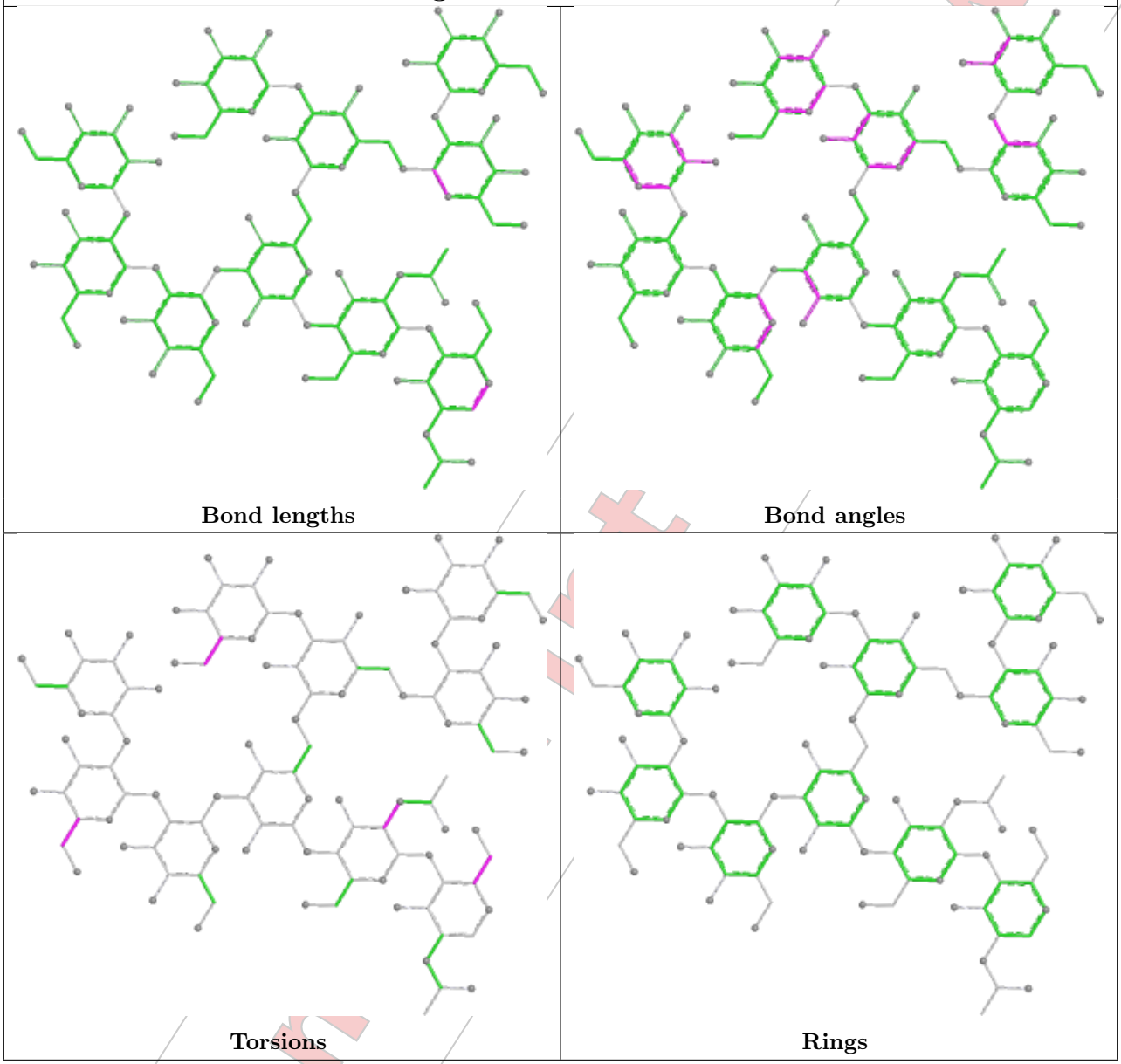


Torsions

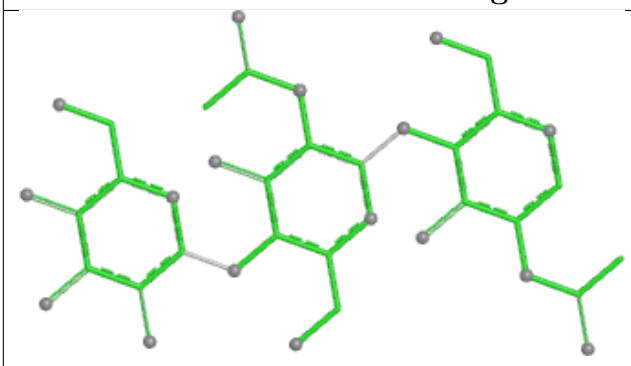


Rings

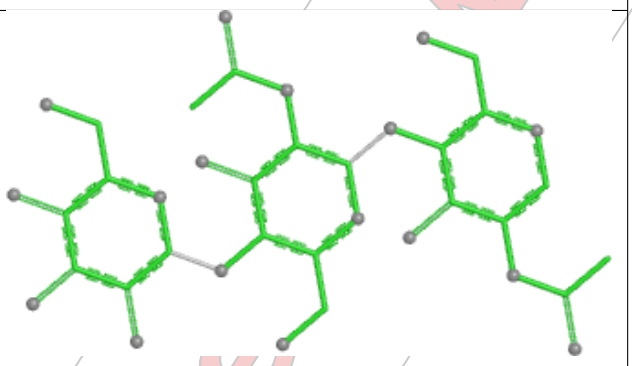
Oligosaccharide Chain G



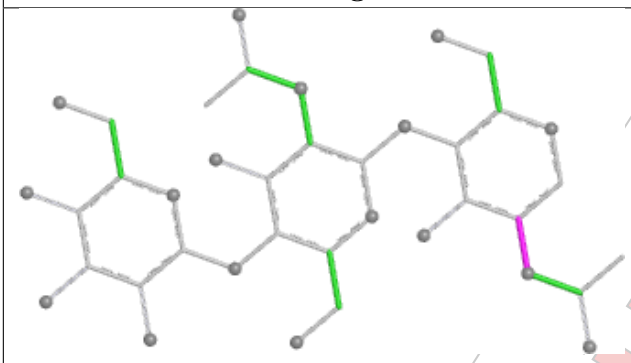
Oligosaccharide Chain M



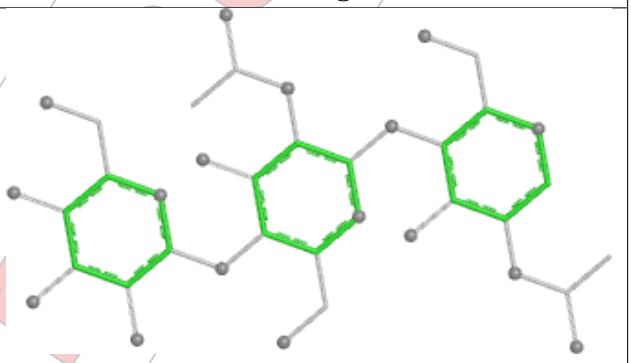
Bond lengths



Bond angles

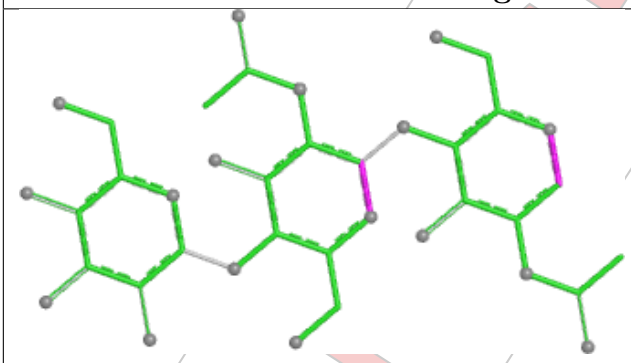


Torsions

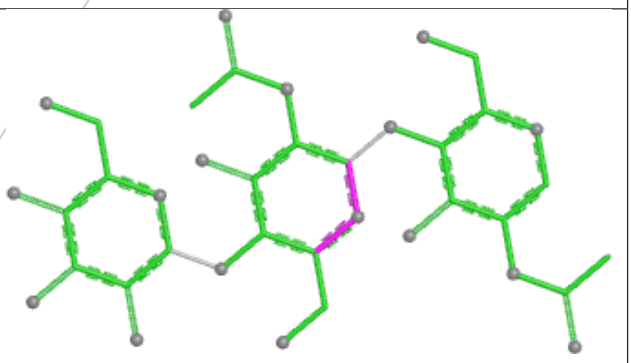


Rings

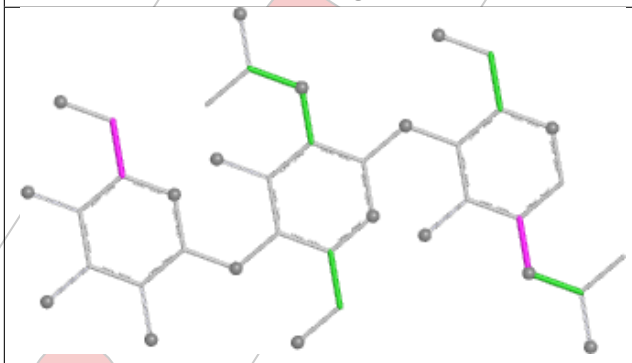
Oligosaccharide Chain H



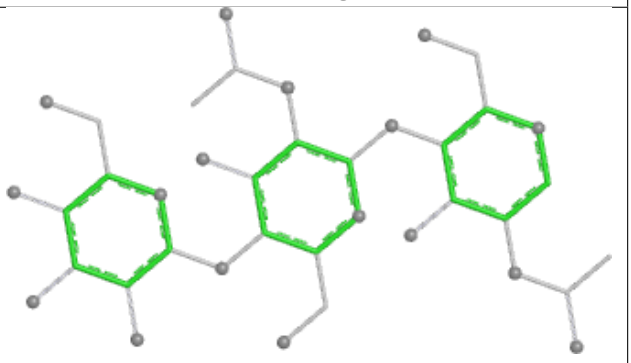
Bond lengths



Bond angles



Torsions



Rings

5.6 Ligand geometry (i)

8 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	D	401	-	6,6,6	0.96	0	6,6,6	1.24	0
11	A1CYU	C	403	-	22,24,24	0.93	1 (4%)	28,33,33	1.75	6 (21%)
10	NAG	E	401	3	14,14,15	0.40	0	17,19,21	0.49	0
9	PTY	B	401	-	40,40,49	2.56	23 (57%)	43,45,54	1.16	3 (6%)
10	NAG	A	402	2	14,14,15	0.18	0	17,19,21	0.48	0
12	ABU	A	401	-	6,6,6	0.83	0	6,6,6	1.41	1 (16%)
9	PTY	C	402	-	40,40,49	2.57	20 (50%)	43,45,54	1.16	3 (6%)
10	NAG	C	401	2	14,14,15	0.29	0	17,19,21	0.42	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
12	ABU	D	401	-	-	2/4/4/4	-
11	A1CYU	C	403	-	-	0/9/25/25	0/3/3/3
10	NAG	E	401	3	-	2/6/23/26	0/1/1/1
9	PTY	B	401	-	-	22/44/44/53	-
10	NAG	A	402	2	-	2/6/23/26	0/1/1/1
12	ABU	A	401	-	-	2/4/4/4	-
9	PTY	C	402	-	-	24/44/44/53	-
10	NAG	C	401	2	-	2/6/23/26	0/1/1/1

All (44) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	C	402	PTY	O7-C8	5.68	1.50	1.34
9	B	401	PTY	O7-C8	5.67	1.50	1.34

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	C	402	PTY	O4-C30	5.51	1.49	1.33
9	B	401	PTY	O4-C30	5.31	1.48	1.33
9	B	401	PTY	C11-C8	4.60	1.64	1.50
9	C	402	PTY	C11-C8	4.46	1.63	1.50
9	C	402	PTY	C31-C30	4.26	1.63	1.50
9	B	401	PTY	C31-C30	4.06	1.62	1.50
11	C	403	A1CYU	C22-C13	-3.78	1.34	1.38
9	C	402	PTY	P1-O14	3.73	1.74	1.59
9	B	401	PTY	P1-O14	3.62	1.73	1.59
9	C	402	PTY	P1-O11	3.51	1.73	1.59
9	B	401	PTY	P1-O11	3.48	1.73	1.59
9	B	401	PTY	C12-C11	3.47	1.64	1.52
9	C	402	PTY	C12-C11	3.46	1.64	1.52
9	C	402	PTY	C1-C6	3.18	1.60	1.50
9	B	401	PTY	C32-C31	3.15	1.63	1.52
9	B	401	PTY	C1-C6	3.14	1.60	1.50
9	C	402	PTY	C2-C3	3.09	1.62	1.49
9	B	401	PTY	C2-C3	3.06	1.62	1.49
9	C	402	PTY	C32-C31	3.05	1.63	1.52
9	C	402	PTY	C5-C6	2.76	1.59	1.50
9	B	401	PTY	C5-C6	2.75	1.59	1.50
9	C	402	PTY	C34-C33	2.60	1.64	1.51
9	B	401	PTY	C35-C34	2.46	1.64	1.51
9	C	402	PTY	C35-C34	2.40	1.63	1.51
9	B	401	PTY	C34-C33	2.39	1.63	1.51
9	B	401	PTY	C13-C12	2.35	1.63	1.51
9	B	401	PTY	C33-C32	2.34	1.63	1.51
9	C	402	PTY	C33-C32	2.32	1.63	1.51
9	C	402	PTY	C18-C17	2.25	1.63	1.51
9	C	402	PTY	C14-C13	2.23	1.62	1.51
9	B	401	PTY	C14-C13	2.23	1.62	1.51
9	B	401	PTY	C18-C17	2.22	1.62	1.51
9	B	401	PTY	C37-C36	2.21	1.62	1.51
9	C	402	PTY	C13-C12	2.18	1.62	1.51
9	B	401	PTY	C39-C38	2.15	1.62	1.51
9	B	401	PTY	C38-C37	2.12	1.62	1.51
9	C	402	PTY	C38-C37	2.10	1.62	1.51
9	C	402	PTY	C39-C38	2.08	1.62	1.51
9	B	401	PTY	C15-C14	2.05	1.62	1.51
9	C	402	PTY	C40-C39	2.05	1.62	1.51
9	B	401	PTY	C17-C16	2.04	1.61	1.51
9	B	401	PTY	C40-C39	2.01	1.61	1.51

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
11	C	403	A1CYU	C10-C2-N3	5.42	126.69	119.77
9	C	402	PTY	O7-C8-C11	3.95	120.02	111.48
9	B	401	PTY	O7-C8-C11	3.93	119.97	111.48
11	C	403	A1CYU	C5-C4-N3	-3.79	104.78	110.65
11	C	403	A1CYU	O1-C2-C10	-3.20	113.14	118.90
11	C	403	A1CYU	C4-N3-C8	-2.91	109.94	114.75
11	C	403	A1CYU	C4-N3-C2	2.60	130.77	123.81
12	A	401	ABU	CB-CG-C	-2.48	108.04	114.51
9	C	402	PTY	O12-P1-O13	-2.46	100.99	112.44
9	B	401	PTY	O12-P1-O13	-2.45	101.04	112.44
9	B	401	PTY	O4-C30-C31	2.42	119.22	111.83
9	C	402	PTY	O4-C30-C31	2.34	118.98	111.83
11	C	403	A1CYU	C9-C8-N3	2.03	114.79	111.55

There are no chirality outliers.

All (56) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
9	B	401	PTY	N1-C2-C3-O11
9	B	401	PTY	C11-C8-O7-C6
9	B	401	PTY	C3-O11-P1-O14
9	C	402	PTY	N1-C2-C3-O11
9	C	402	PTY	C11-C8-O7-C6
9	C	402	PTY	O30-C30-O4-C1
9	C	402	PTY	O10-C8-O7-C6
9	C	402	PTY	C31-C30-O4-C1
9	B	401	PTY	O10-C8-O7-C6
10	A	402	NAG	C4-C5-C6-O6
9	B	401	PTY	C31-C30-O4-C1
9	B	401	PTY	O30-C30-O4-C1
10	A	402	NAG	O5-C5-C6-O6
10	E	401	NAG	C8-C7-N2-C2
10	E	401	NAG	O7-C7-N2-C2
10	C	401	NAG	O5-C5-C6-O6
9	C	402	PTY	C31-C32-C33-C34
9	B	401	PTY	C14-C15-C16-C17
9	B	401	PTY	C34-C35-C36-C37
9	B	401	PTY	C12-C13-C14-C15
9	B	401	PTY	C8-C11-C12-C13
9	B	401	PTY	C36-C37-C38-C39
9	C	402	PTY	C14-C15-C16-C17

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Mol	Chain	Res	Type	Atoms
9	B	401	PTY	C15-C16-C17-C18
9	C	402	PTY	C40-C41-C42-C43
9	C	402	PTY	C12-C13-C14-C15
9	B	401	PTY	C38-C39-C40-C41
10	C	401	NAG	C4-C5-C6-O6
9	C	402	PTY	C34-C35-C36-C37
9	B	401	PTY	C37-C38-C39-C40
9	B	401	PTY	C30-C31-C32-C33
9	C	402	PTY	C15-C16-C17-C18
9	B	401	PTY	C13-C14-C15-C16
9	B	401	PTY	C2-C3-O11-P1
9	C	402	PTY	C2-C3-O11-P1
9	C	402	PTY	C32-C33-C34-C35
9	B	401	PTY	C17-C18-C19-C20
9	C	402	PTY	C16-C17-C18-C19
9	C	402	PTY	C36-C37-C38-C39
9	C	402	PTY	C41-C42-C43-C44
9	B	401	PTY	C3-O11-P1-O13
9	C	402	PTY	C3-O11-P1-O12
9	C	402	PTY	C3-O11-P1-O13
9	C	402	PTY	C3-O11-P1-O14
9	C	402	PTY	C39-C40-C41-C42
9	C	402	PTY	C38-C39-C40-C41
9	B	401	PTY	C35-C36-C37-C38
9	B	401	PTY	C31-C32-C33-C34
12	D	401	ABU	OXT-C-CG-CB
12	D	401	ABU	O-C-CG-CB
12	A	401	ABU	O-C-CG-CB
9	C	402	PTY	O4-C1-C6-O7
12	A	401	ABU	OXT-C-CG-CB
9	C	402	PTY	O4-C1-C6-C5
9	B	401	PTY	C33-C34-C35-C36
9	C	402	PTY	C30-C31-C32-C33

There are no ring outliers.

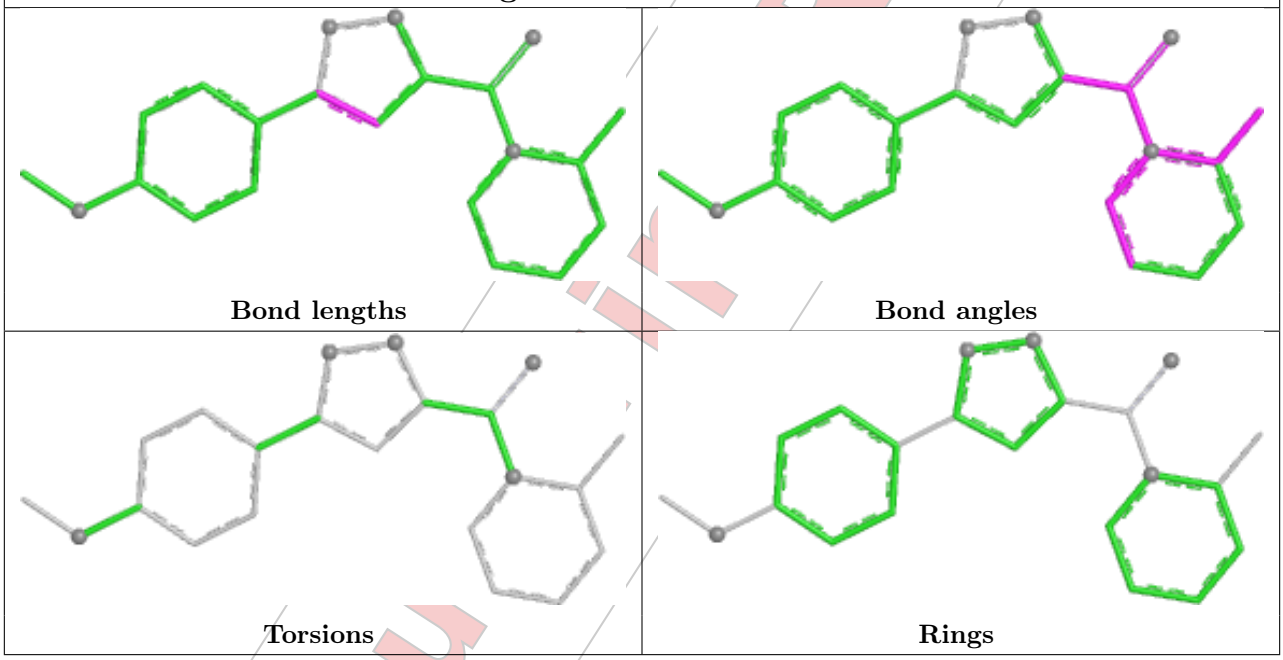
2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
10	A	402	NAG	1	0
12	A	401	ABU	1	0

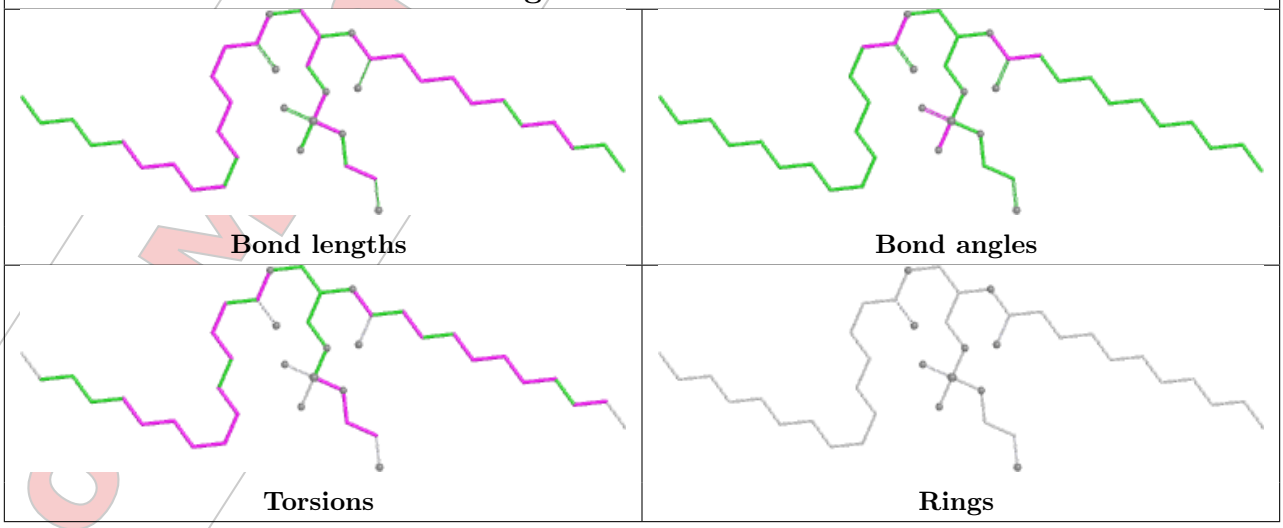
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths,

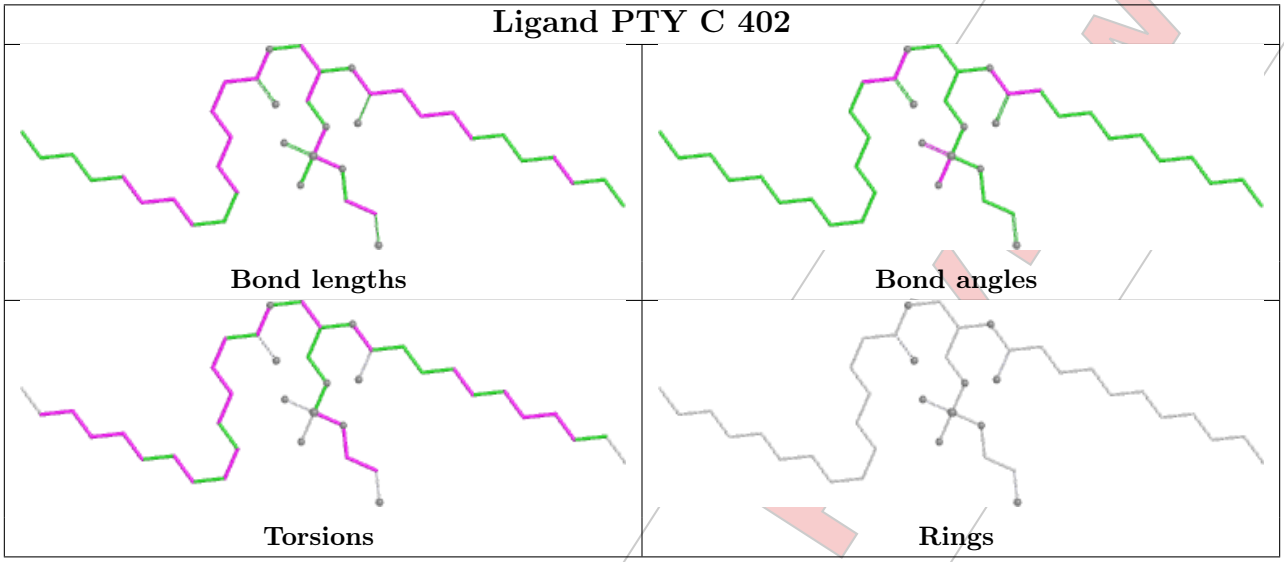
bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less then 5% of the Mogul distribution of torsion angles is within 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 A1CYU C 403



Ligand PTY B 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.

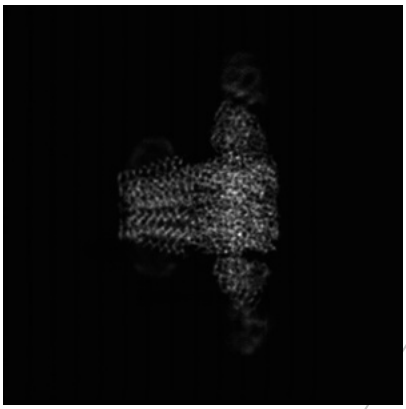
6 Map visualisation

This section contains visualisations of the EMDB entry EMD-73211. These allow visual inspection of the internal detail of the map and identification of artifacts.

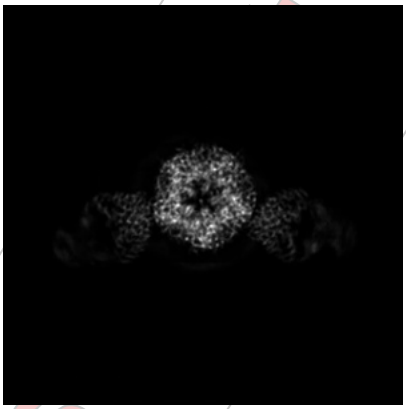
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections

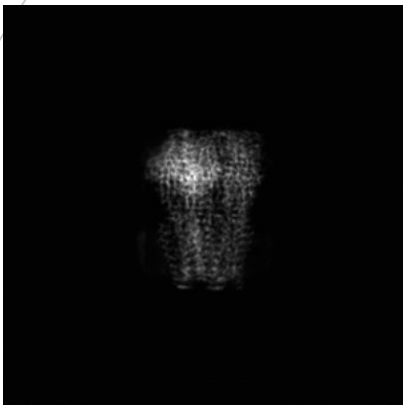
6.1.1 Primary map



X

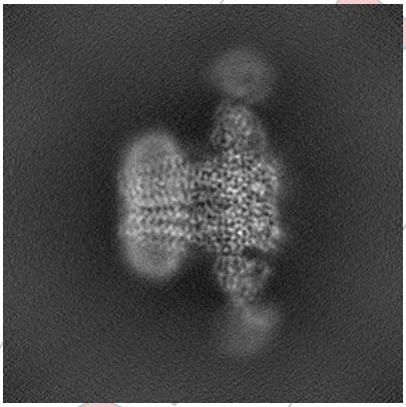


Y

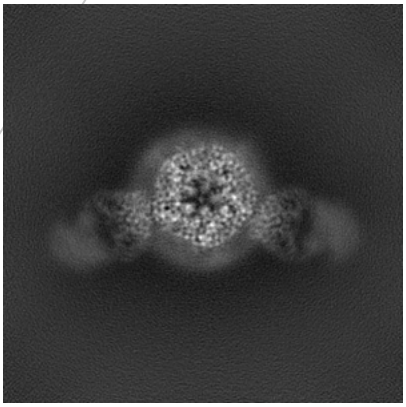


Z

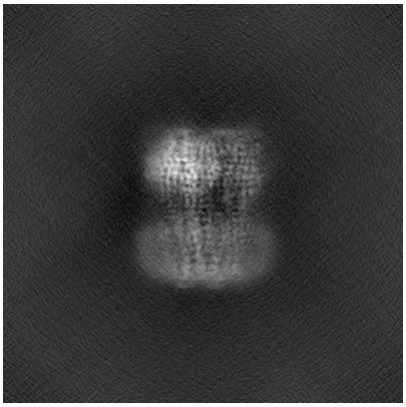
6.1.2 Raw map



X



Y



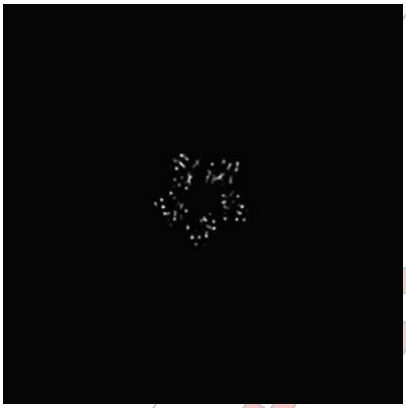
Z

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

6.2 Central slices (i)

6.2.1 Primary map

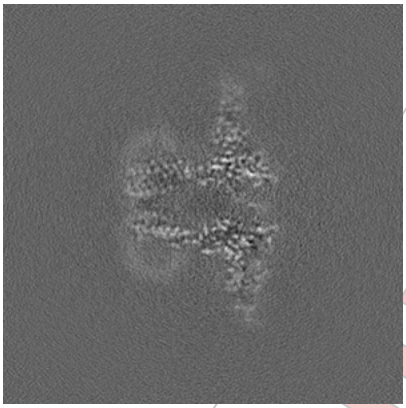
X Index: 144



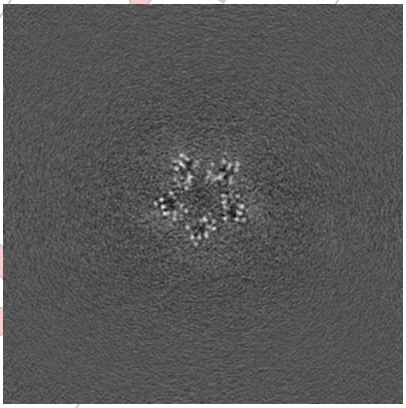
Y Index: 144

Z Index: 144

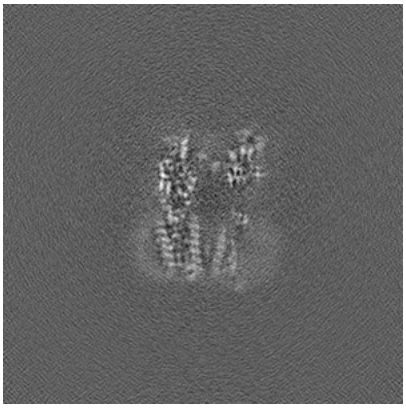
6.2.2 Raw map



X Index: 144



Y Index: 144



Z Index: 144

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

6.3 Largest variance slices (i)

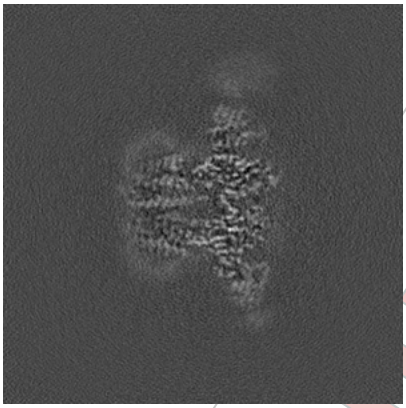
6.3.1 Primary map

X Index: 137

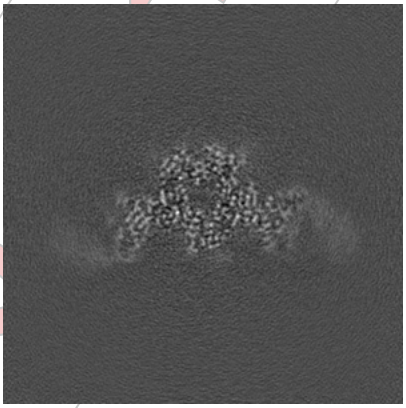
Y Index: 163

Z Index: 125

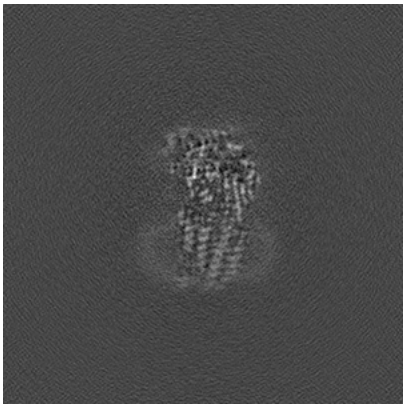
6.3.2 Raw map



X Index: 137



Y Index: 163

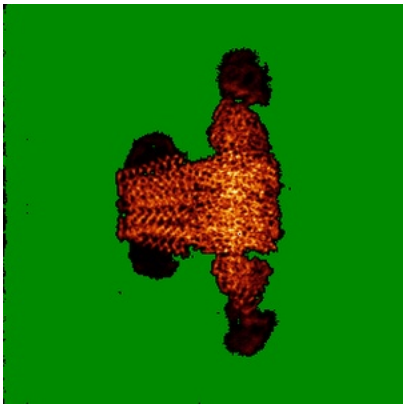


Z Index: 125

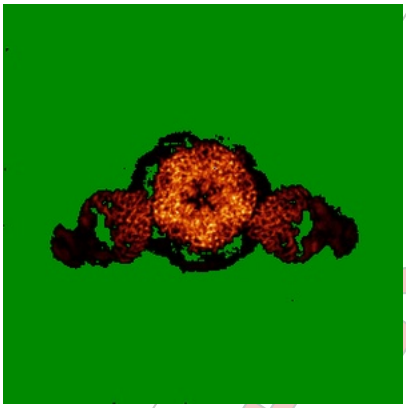
The images above show the largest variance slices of the map in three orthogonal directions.

6.4 Orthogonal standard-deviation projections (False-color)

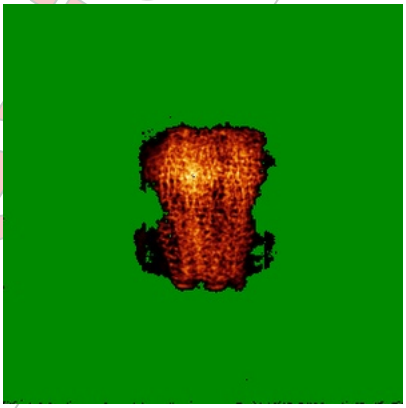
6.4.1 Primary map



X

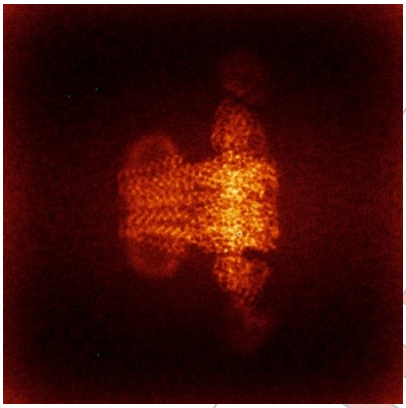


Y

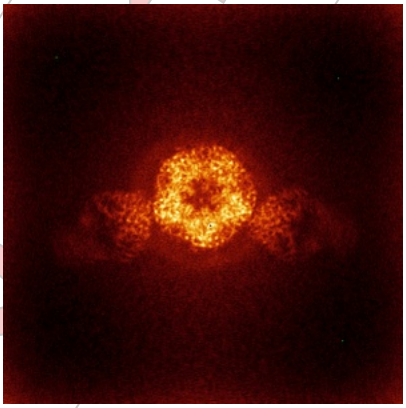


Z

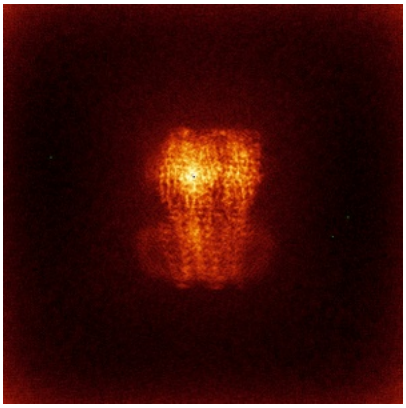
6.4.2 Raw map



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.

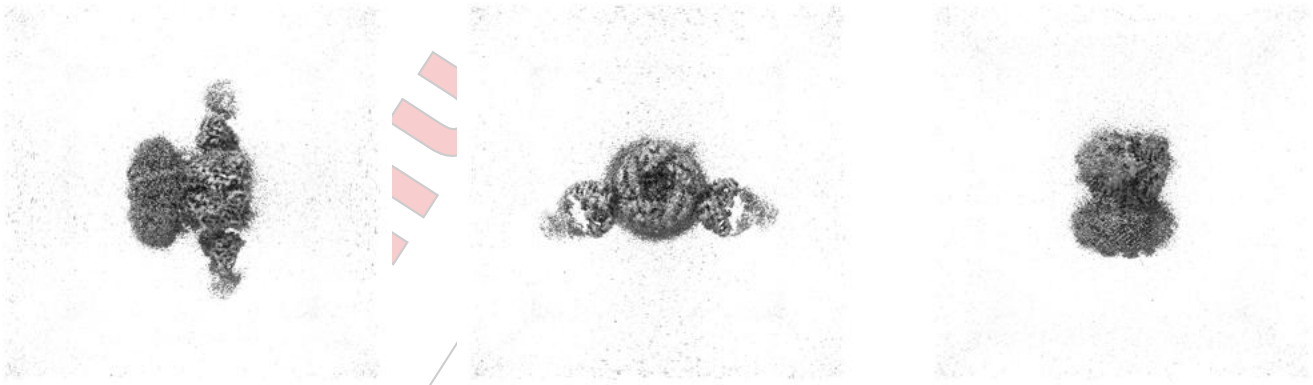
6.5 Orthogonal surface views (i)

6.5.1 Primary map

X Y Z

The images above show the 3D surface view of the map at the recommended contour level 0.0297. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.5.2 Raw map



X Y Z

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

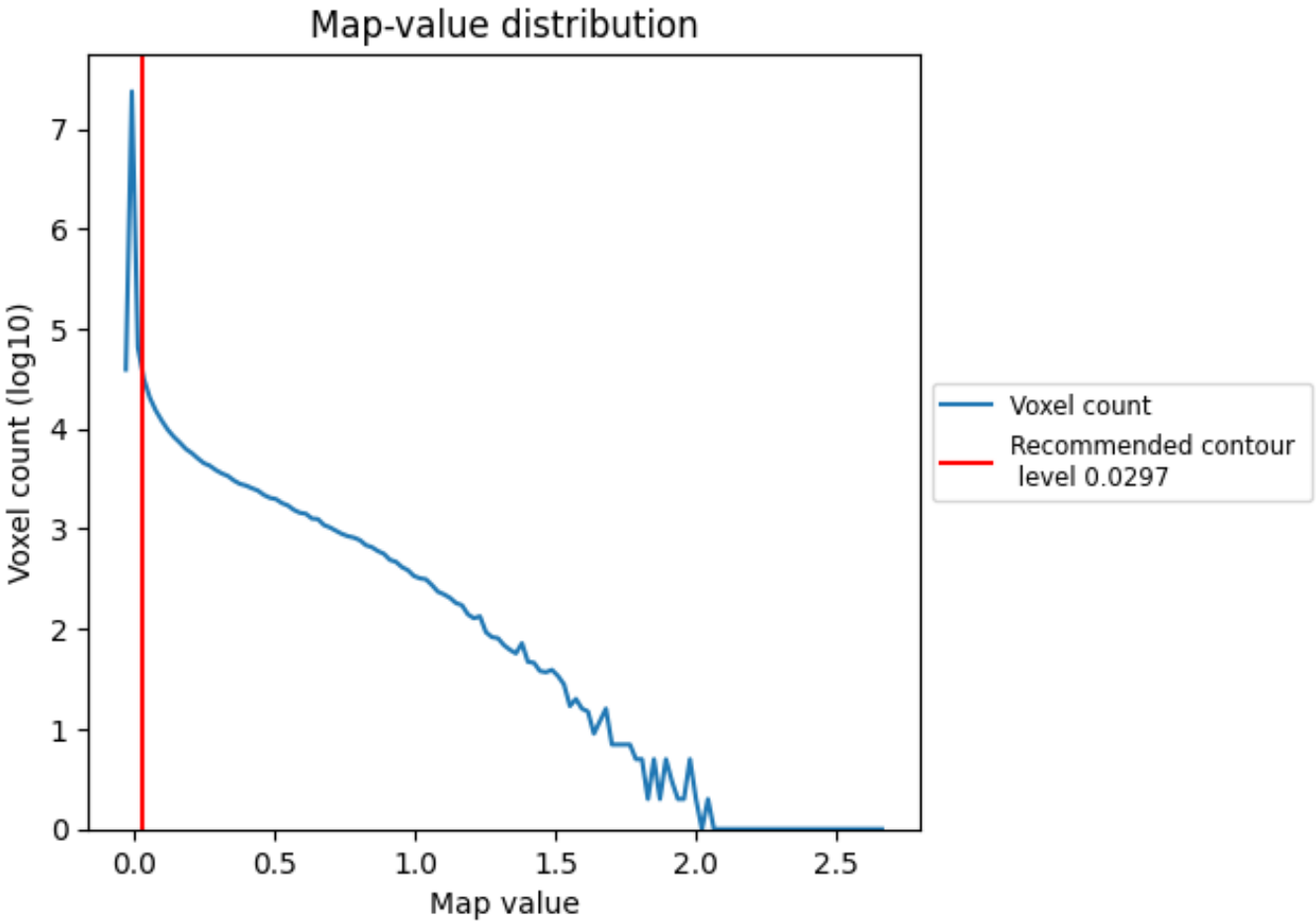
6.6 Mask visualisation (i)

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

7 Map analysis

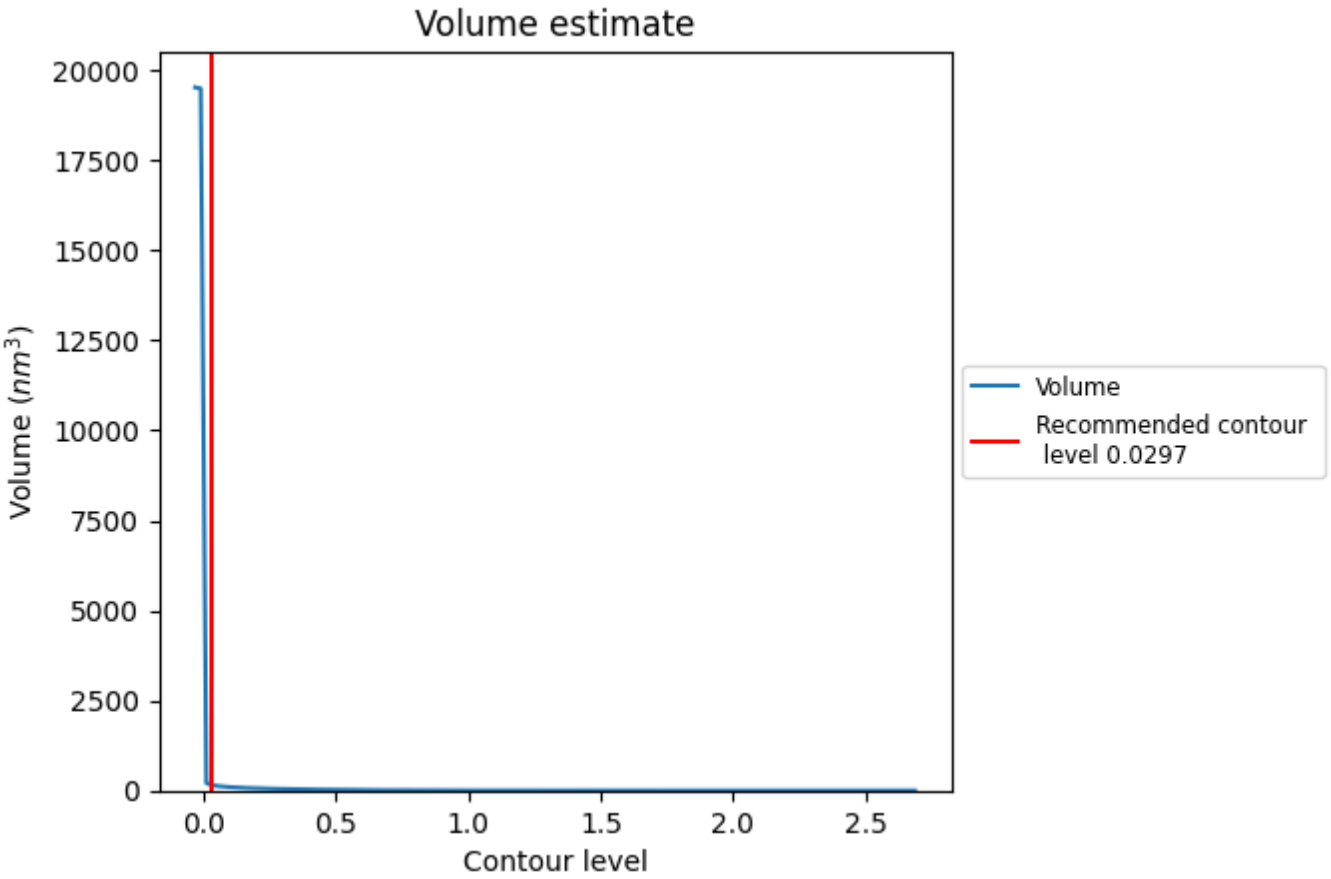
This section contains the results of statistical analysis of the map.

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.

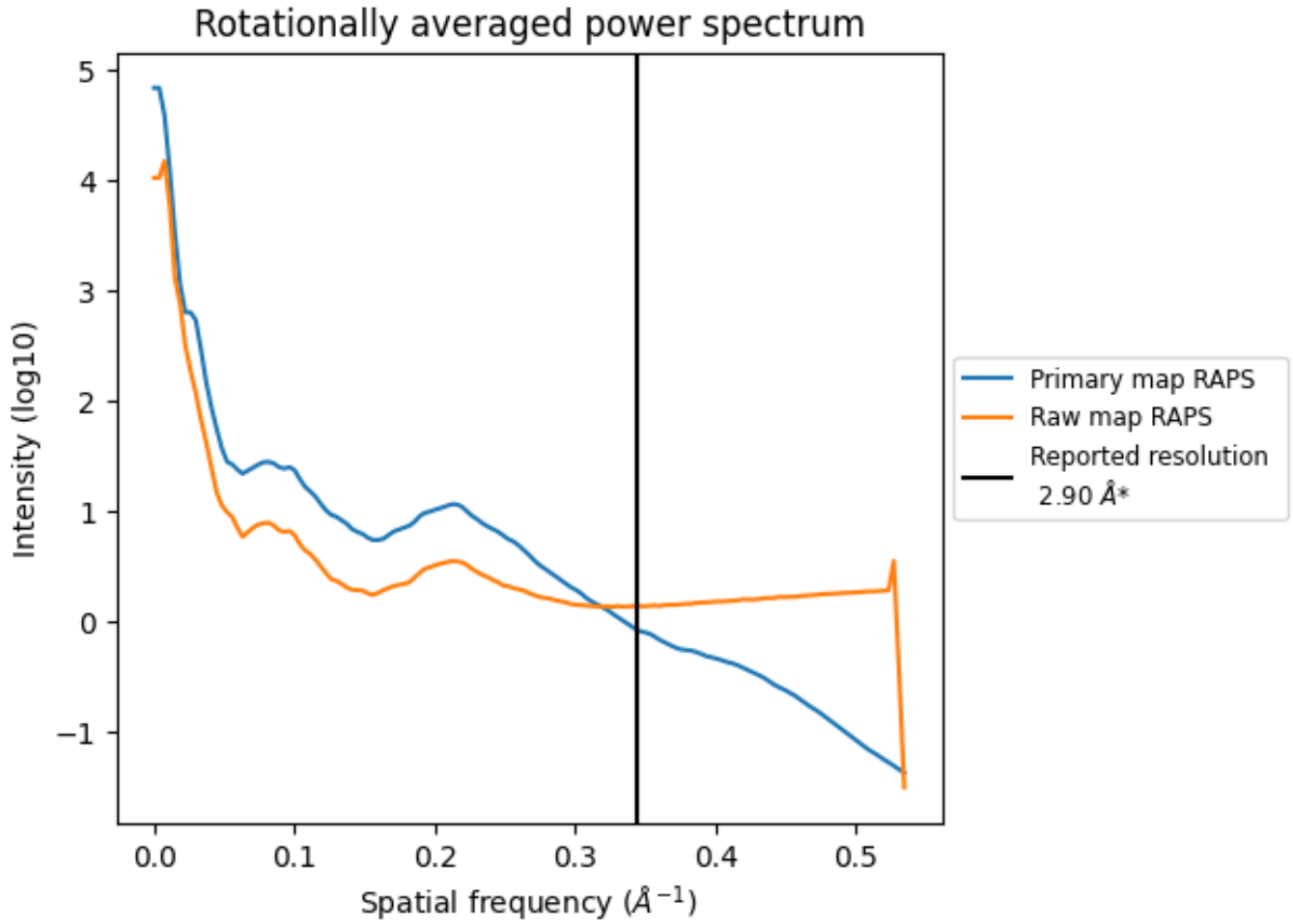
7.2 Volume estimate (i)



The volume at the recommended contour level is 162 nm^3; this corresponds to an approximate mass of 146 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)

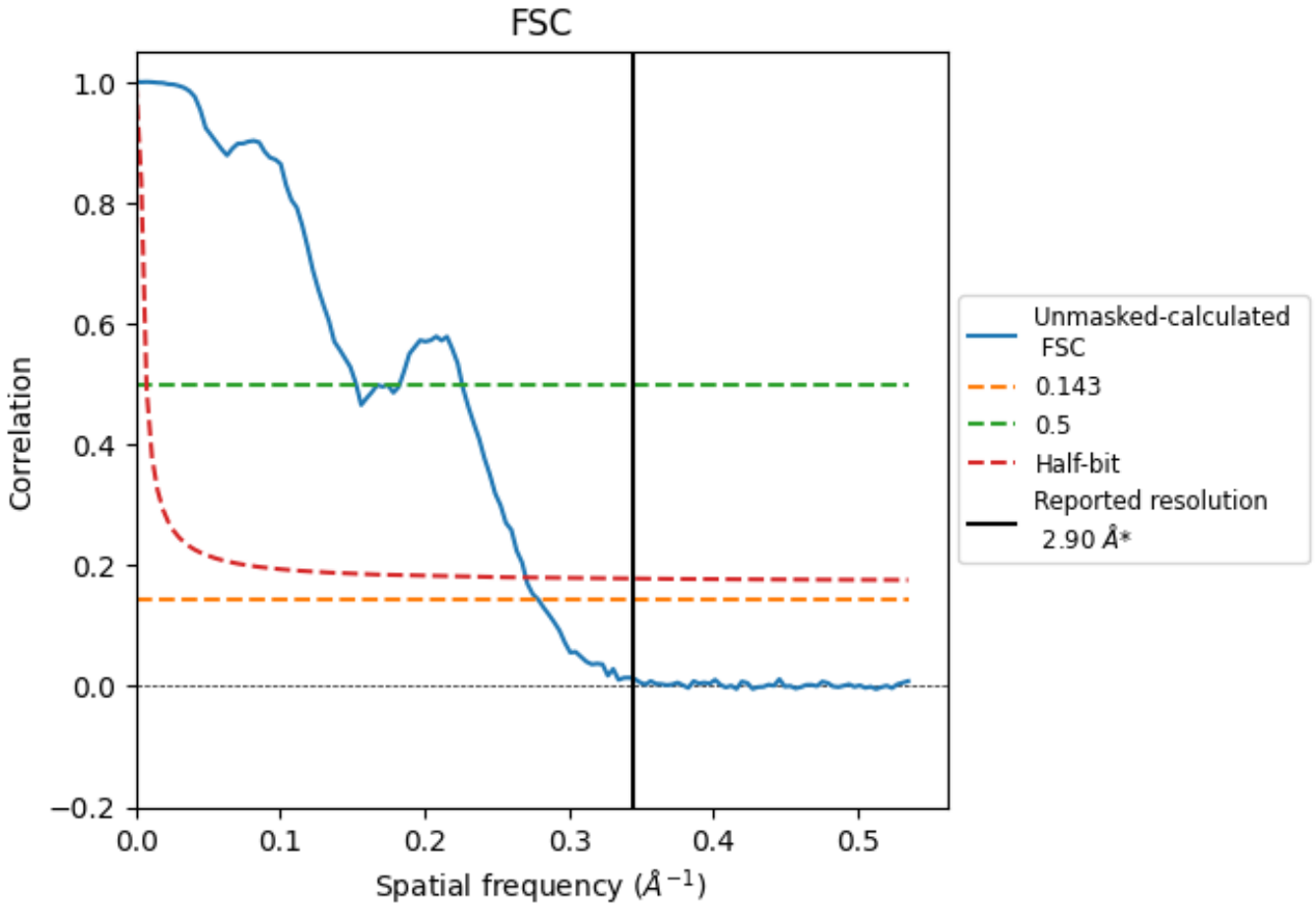


*Reported resolution corresponds to spatial frequency of 0.345 Å^{-1}

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 0.345 Å^{-1}

8.2 Resolution estimates (i)

Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.90	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	3.59	6.56	3.70

*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 3.59 differs from the reported value 2.9 by more than 10 %

9 Map-model fit

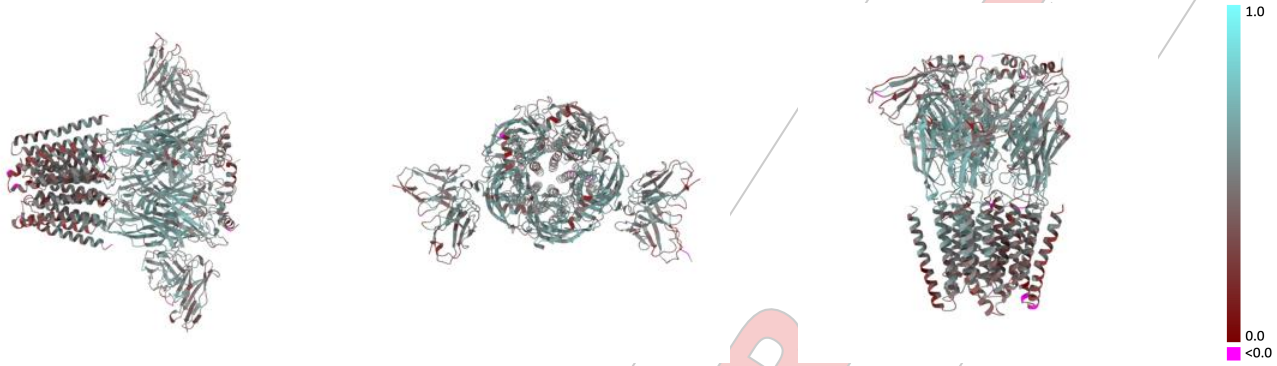
This section contains information regarding the fit between EMDB map EMD-73211 and PDB model 9YNN. Per-residue inclusion information can be found in section 3 on page 12.

9.1 Map-model overlay

X Y Z

The images above show the 3D surface view of the map at the recommended contour level 0.0297 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.

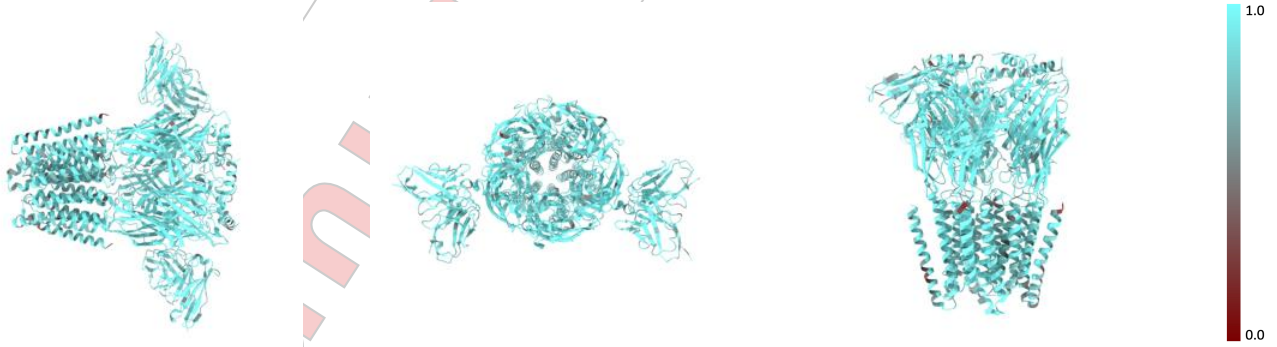
9.2 Q-score mapped to coordinate model



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.

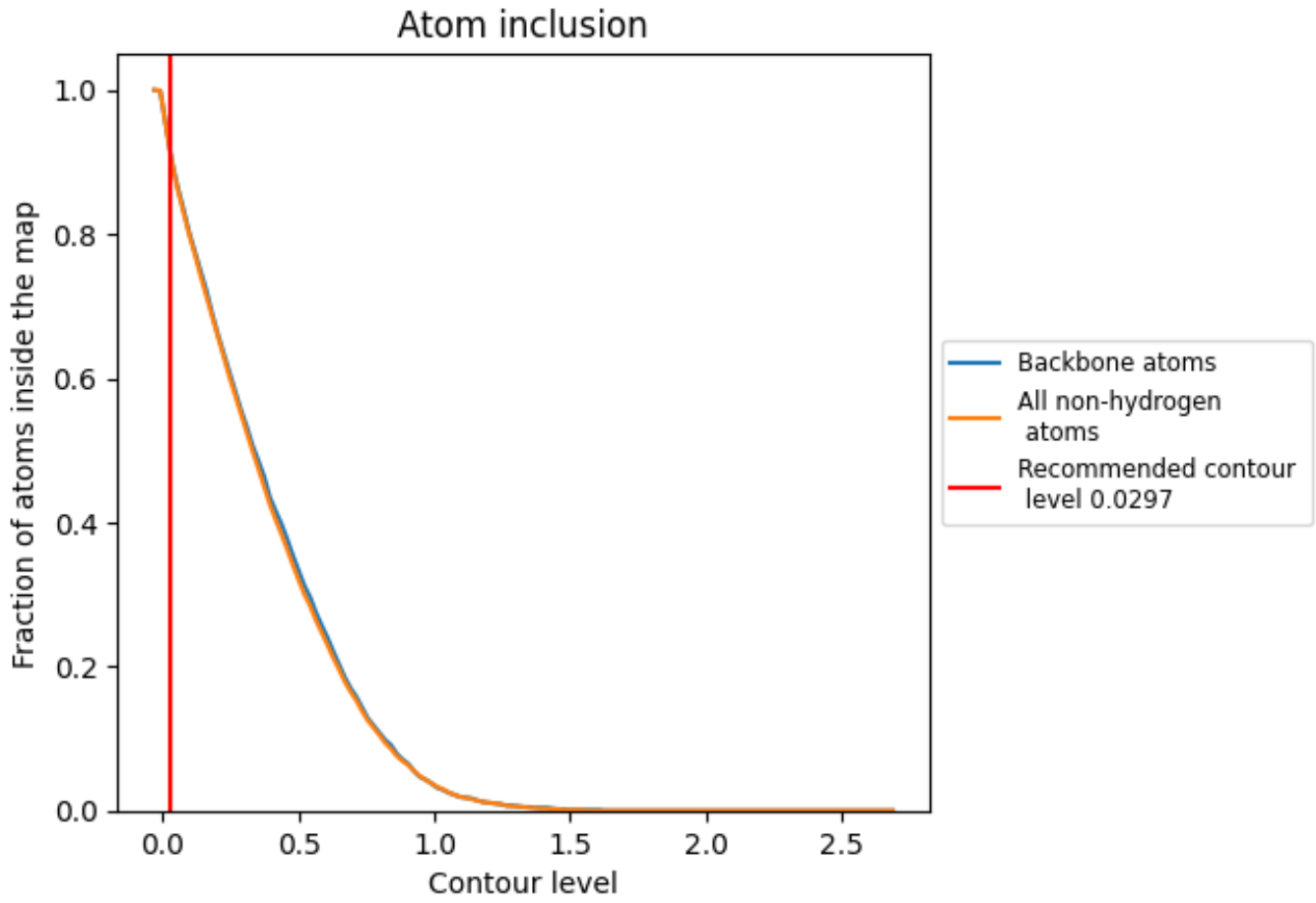
9.3 Atom inclusion mapped to coordinate model



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 (0.0297).

9.4 Atom inclusion



At the recommended contour level, 91% of all backbone atoms, 91% 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 (0.0297) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.9080	0.4930
A	0.9170	0.4880
B	0.9230	0.5180
C	0.9150	0.5160
D	0.9300	0.5280
E	0.8870	0.4580
F	0.8930	0.5020
G	0.8190	0.4230
H	0.7690	0.4200
I	0.8960	0.4800
J	0.8870	0.4560
K	0.8630	0.4390
L	0.8990	0.4700
M	0.8720	0.5160