



Full wwPDB X-ray Structure Validation Report ⓘ

Jul 21, 2025 – 01:17 PM JST

PDB ID : 9IWU / pdb_00009iwu
Title : CTB10-PE3.0-(R)-1g complex
Authors : Fu, K.; Rao, Y.J.
Deposited on : 2024-07-26
Resolution : 2.20 Å(reported)

This is a Full wwPDB X-ray Structure 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/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

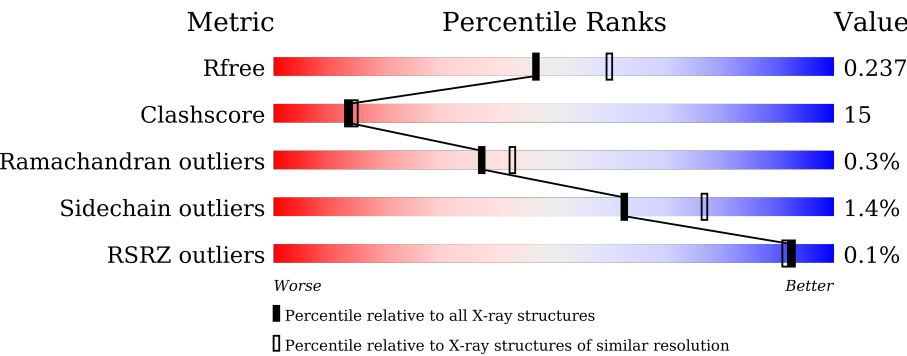
MolProbity : 4-5-2 with Phenix2.0rc1
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 2.0rc1
EDS : 3.0
buster-report : 1.1.7 (2018)
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4 : 9.0.006 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.44

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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)	Similar resolution (#Entries, resolution range(Å))
R_{free}	164625	5791 (2.20-2.20)
Clashscore	180529	6634 (2.20-2.20)
Ramachandran outliers	177936	6560 (2.20-2.20)
Sidechain outliers	177891	6561 (2.20-2.20)
RSRZ outliers	164620	5791 (2.20-2.20)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	141	<div><div></div><div></div><div></div><div></div><div></div></div> <div>64%24%12%</div>
1	B	141	<div><div></div><div></div><div></div><div></div><div></div></div> <div>69%26%6%</div>
1	C	141	<div><div></div><div></div><div></div><div></div><div></div></div> <div>64%23%•11%</div>
1	D	141	<div><div></div><div></div><div></div><div></div><div></div></div> <div>67%23%•10%</div>
1	E	141	<div><div></div><div></div><div></div><div></div><div></div></div> <div>57%30%•12%</div>
1	F	141	<div><div></div><div></div><div></div><div></div><div></div></div> <div>66%21%•11%</div>

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Mol	Chain	Length	Quality of chain
1	G	141	<div><div>%</div><div><div></div><div>69%</div><div>18%</div><div>•</div><div>12%</div></div></div>
1	H	141	<div><div></div><div>60%</div><div>26%</div><div>•</div><div>12%</div></div>

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 8857 atoms, of which 142 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, 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 CTB10.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	124	Total	C	N	O	S	0	0	0
			1018	658	165	190	5			
1	B	133	Total	C	N	O	S	0	0	0
			1088	700	178	205	5			
1	C	125	Total	C	N	O	S	0	0	0
			1032	666	171	190	5			
1	D	127	Total	C	N	O	S	0	0	0
			1046	676	172	193	5			
1	E	124	Total	C	N	O	S	0	0	0
			1029	665	169	190	5			
1	F	125	Total	C	N	O	S	0	0	0
			1033	667	171	190	5			
1	G	124	Total	C	N	O	S	0	0	0
			1025	663	169	188	5			
1	H	124	Total	C	N	O	S	0	0	0
			1016	657	168	186	5			

There are 88 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	40	PBF	MET	engineered mutation	UNP A0A977K7H6
A	86	ILE	MET	engineered mutation	UNP A0A977K7H6
A	133	ALA	-	expression tag	UNP A0A977K7H6
A	134	LEU	-	expression tag	UNP A0A977K7H6
A	135	GLU	-	expression tag	UNP A0A977K7H6
A	136	HIS	-	expression tag	UNP A0A977K7H6
A	137	HIS	-	expression tag	UNP A0A977K7H6
A	138	HIS	-	expression tag	UNP A0A977K7H6
A	139	HIS	-	expression tag	UNP A0A977K7H6
A	140	HIS	-	expression tag	UNP A0A977K7H6
A	141	HIS	-	expression tag	UNP A0A977K7H6
B	40	PBF	MET	engineered mutation	UNP A0A977K7H6
B	86	ILE	MET	engineered mutation	UNP A0A977K7H6

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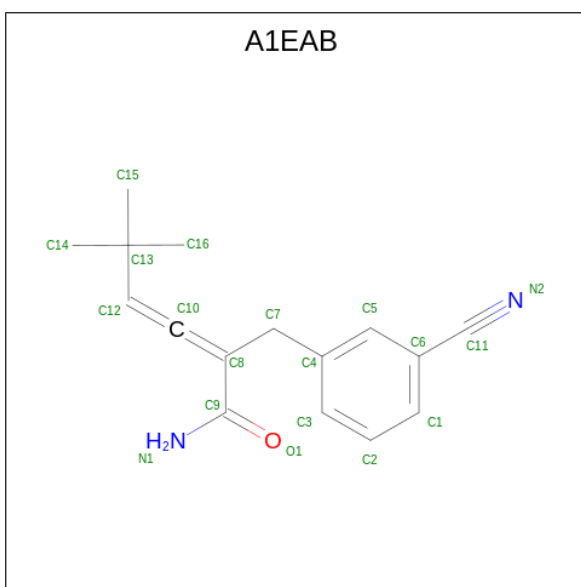
Chain	Residue	Modelled	Actual	Comment	Reference
B	133	ALA	-	expression tag	UNP A0A977K7H6
B	134	LEU	-	expression tag	UNP A0A977K7H6
B	135	GLU	-	expression tag	UNP A0A977K7H6
B	136	HIS	-	expression tag	UNP A0A977K7H6
B	137	HIS	-	expression tag	UNP A0A977K7H6
B	138	HIS	-	expression tag	UNP A0A977K7H6
B	139	HIS	-	expression tag	UNP A0A977K7H6
B	140	HIS	-	expression tag	UNP A0A977K7H6
B	141	HIS	-	expression tag	UNP A0A977K7H6
C	40	PBF	MET	engineered mutation	UNP A0A977K7H6
C	86	ILE	MET	engineered mutation	UNP A0A977K7H6
C	133	ALA	-	expression tag	UNP A0A977K7H6
C	134	LEU	-	expression tag	UNP A0A977K7H6
C	135	GLU	-	expression tag	UNP A0A977K7H6
C	136	HIS	-	expression tag	UNP A0A977K7H6
C	137	HIS	-	expression tag	UNP A0A977K7H6
C	138	HIS	-	expression tag	UNP A0A977K7H6
C	139	HIS	-	expression tag	UNP A0A977K7H6
C	140	HIS	-	expression tag	UNP A0A977K7H6
C	141	HIS	-	expression tag	UNP A0A977K7H6
D	40	PBF	MET	engineered mutation	UNP A0A977K7H6
D	86	ILE	MET	engineered mutation	UNP A0A977K7H6
D	133	ALA	-	expression tag	UNP A0A977K7H6
D	134	LEU	-	expression tag	UNP A0A977K7H6
D	135	GLU	-	expression tag	UNP A0A977K7H6
D	136	HIS	-	expression tag	UNP A0A977K7H6
D	137	HIS	-	expression tag	UNP A0A977K7H6
D	138	HIS	-	expression tag	UNP A0A977K7H6
D	139	HIS	-	expression tag	UNP A0A977K7H6
D	140	HIS	-	expression tag	UNP A0A977K7H6
D	141	HIS	-	expression tag	UNP A0A977K7H6
E	40	PBF	MET	engineered mutation	UNP A0A977K7H6
E	86	ILE	MET	engineered mutation	UNP A0A977K7H6
E	133	ALA	-	expression tag	UNP A0A977K7H6
E	134	LEU	-	expression tag	UNP A0A977K7H6
E	135	GLU	-	expression tag	UNP A0A977K7H6
E	136	HIS	-	expression tag	UNP A0A977K7H6
E	137	HIS	-	expression tag	UNP A0A977K7H6
E	138	HIS	-	expression tag	UNP A0A977K7H6
E	139	HIS	-	expression tag	UNP A0A977K7H6
E	140	HIS	-	expression tag	UNP A0A977K7H6
E	141	HIS	-	expression tag	UNP A0A977K7H6

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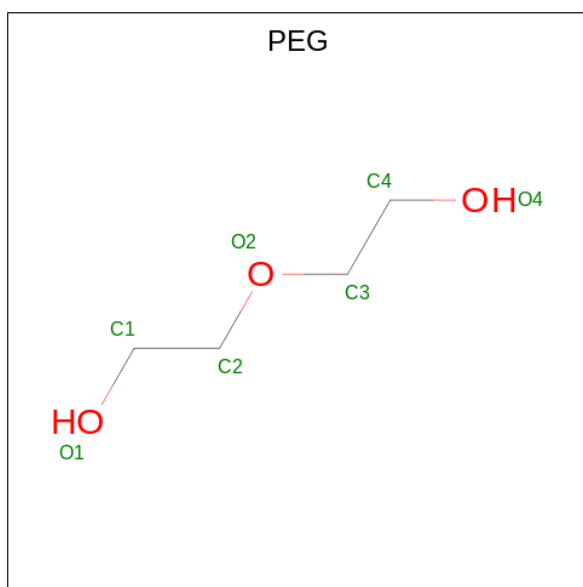
Chain	Residue	Modelled	Actual	Comment	Reference
F	40	PBF	MET	engineered mutation	UNP A0A977K7H6
F	86	ILE	MET	engineered mutation	UNP A0A977K7H6
F	133	ALA	-	expression tag	UNP A0A977K7H6
F	134	LEU	-	expression tag	UNP A0A977K7H6
F	135	GLU	-	expression tag	UNP A0A977K7H6
F	136	HIS	-	expression tag	UNP A0A977K7H6
F	137	HIS	-	expression tag	UNP A0A977K7H6
F	138	HIS	-	expression tag	UNP A0A977K7H6
F	139	HIS	-	expression tag	UNP A0A977K7H6
F	140	HIS	-	expression tag	UNP A0A977K7H6
F	141	HIS	-	expression tag	UNP A0A977K7H6
G	40	PBF	MET	engineered mutation	UNP A0A977K7H6
G	86	ILE	MET	engineered mutation	UNP A0A977K7H6
G	133	ALA	-	expression tag	UNP A0A977K7H6
G	134	LEU	-	expression tag	UNP A0A977K7H6
G	135	GLU	-	expression tag	UNP A0A977K7H6
G	136	HIS	-	expression tag	UNP A0A977K7H6
G	137	HIS	-	expression tag	UNP A0A977K7H6
G	138	HIS	-	expression tag	UNP A0A977K7H6
G	139	HIS	-	expression tag	UNP A0A977K7H6
G	140	HIS	-	expression tag	UNP A0A977K7H6
G	141	HIS	-	expression tag	UNP A0A977K7H6
H	40	PBF	MET	engineered mutation	UNP A0A977K7H6
H	86	ILE	MET	engineered mutation	UNP A0A977K7H6
H	133	ALA	-	expression tag	UNP A0A977K7H6
H	134	LEU	-	expression tag	UNP A0A977K7H6
H	135	GLU	-	expression tag	UNP A0A977K7H6
H	136	HIS	-	expression tag	UNP A0A977K7H6
H	137	HIS	-	expression tag	UNP A0A977K7H6
H	138	HIS	-	expression tag	UNP A0A977K7H6
H	139	HIS	-	expression tag	UNP A0A977K7H6
H	140	HIS	-	expression tag	UNP A0A977K7H6
H	141	HIS	-	expression tag	UNP A0A977K7H6

- Molecule 2 is 2-[(3-cyanophenyl)methyl]-5,5-dimethyl-hexa-2,3-dienamide (CCD ID: A1EAB) (formula: C₁₆H₁₈N₂O) (labeled as "Ligand of Interest" by depositor).



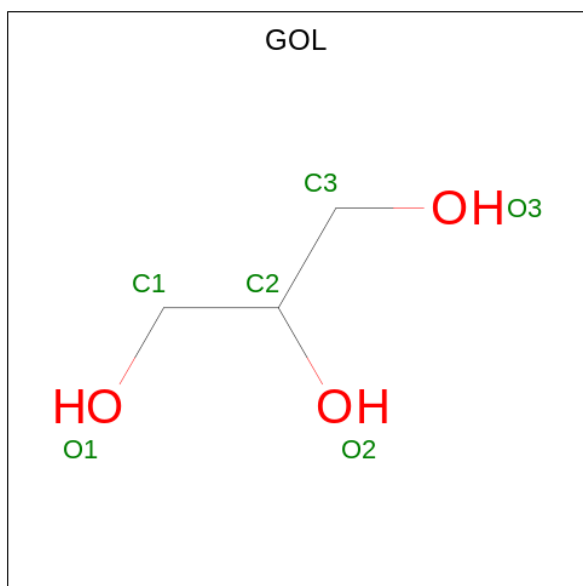
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	H	N	O	0	0
			37	16	18	2	1		
2	C	1	Total	C	H	N	O	0	0
			37	16	18	2	1		
2	E	1	Total	C	H	N	O	0	0
			37	16	18	2	1		
2	F	1	Total	C	H	N	O	0	0
			37	16	18	2	1		
2	G	1	Total	C	H	N	O	0	0
			37	16	18	2	1		
2	H	1	Total	C	H	N	O	0	0
			37	16	18	2	1		

- Molecule 3 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: $C_4H_{10}O_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	B	1	Total	C	H	O	0	0
			17	4	10	3		

- Molecule 4 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	F	1	Total	C	H	O	0	0
			14	3	8	3		
4	H	1	Total	C	H	O	0	0
			14	3	8	3		
4	H	1	Total	C	H	O	0	0
			14	3	8	3		

- Molecule 5 is water.

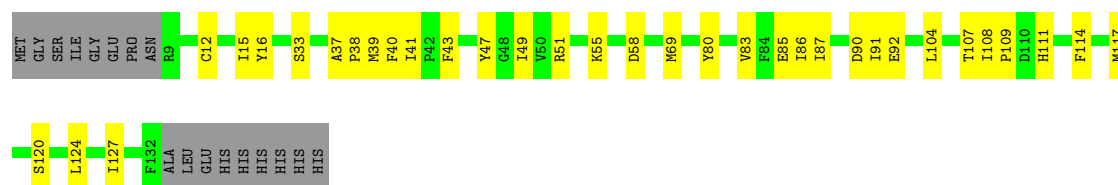
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	36	Total 36	O 36	0	0
5	B	40	Total 40	O 40	0	0
5	C	38	Total 38	O 38	0	0
5	D	41	Total 41	O 41	0	0
5	E	32	Total 32	O 32	0	0
5	F	35	Total 35	O 35	0	0
5	G	23	Total 23	O 23	0	0
5	H	44	Total 44	O 44	0	0

3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron 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 dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). 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.

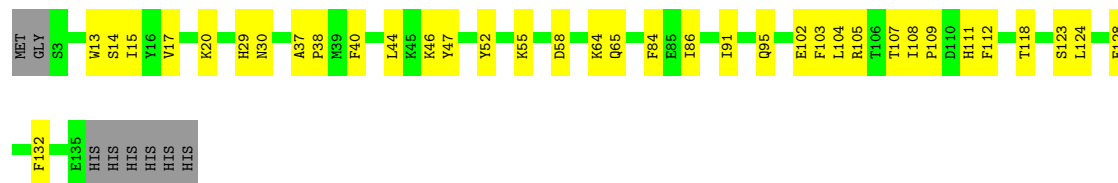
• Molecule 1: CTB10

Chain A: 



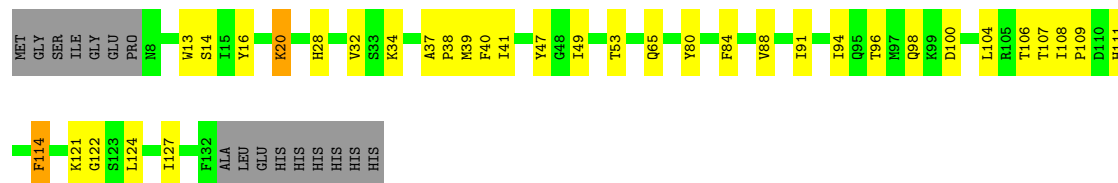
• Molecule 1: CTB10

Chain B: 



• Molecule 1: CTB10

Chain C: 



• Molecule 1: CTB10

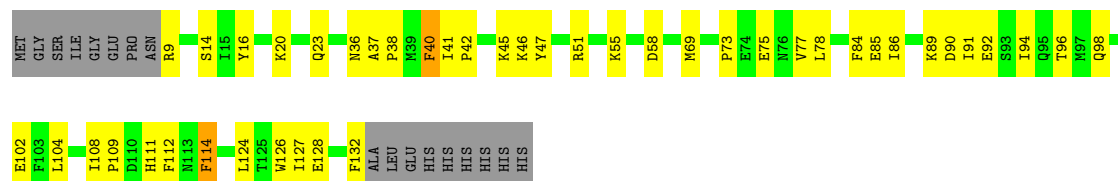
Chain D: 





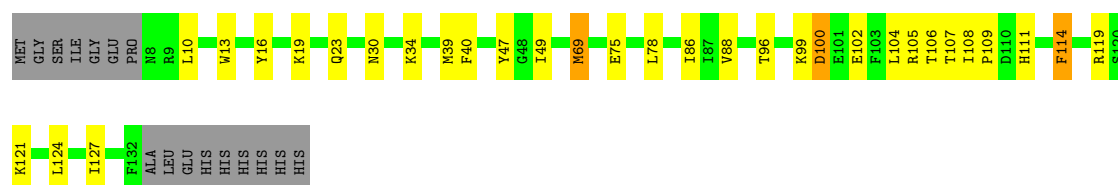
• Molecule 1: CTB10

Chain E: 57% 30% 12%



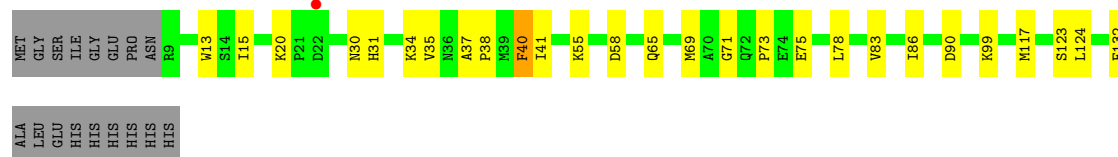
• Molecule 1: CTB10

Chain F: 66% 21% 11%



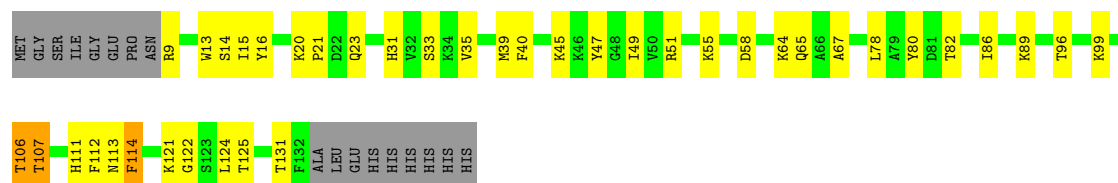
• Molecule 1: CTB10

Chain G: 69% 18% 12%



• Molecule 1: CTB10

Chain H: 60% 26% 12%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	39.13Å 163.71Å 87.41Å 90.00° 93.01° 90.00°	Depositor
Resolution (Å)	46.27 – 2.20 46.27 – 2.20	Depositor EDS
% Data completeness (in resolution range)	97.7 (46.27-2.20) 97.8 (46.27-2.20)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.82 (at 2.20Å)	Xtriage
Refinement program	PHENIX (1.20.1_4487: ???)	Depositor
R, R_{free}	0.182 , 0.238 0.189 , 0.237	Depositor DCC
R_{free} test set	2845 reflections (5.11%)	wwPDB-VP
Wilson B-factor (Å ²)	30.0	Xtriage
Anisotropy	0.569	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 36.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	8857	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 13.20% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, PEG, A1EAB, PBF

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.69	0/1024	0.83	0/1387
1	B	0.74	2/1095 (0.2%)	0.87	0/1483
1	C	0.66	0/1038	0.82	0/1404
1	D	0.71	0/1052	0.86	0/1423
1	E	0.72	0/1035	0.86	0/1399
1	F	0.78	0/1039	1.06	1/1405 (0.1%)
1	G	0.67	0/1031	0.85	0/1394
1	H	0.79	2/1022 (0.2%)	0.98	3/1384 (0.2%)
All	All	0.72	4/8336 (0.0%)	0.90	4/11279 (0.0%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	29	HIS	C-N	-6.99	1.24	1.33
1	H	67	ALA	C-N	-6.78	1.24	1.33
1	H	13	TRP	C-N	-5.86	1.25	1.33
1	B	30	ASN	C-N	5.60	1.41	1.33

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	106	THR	CA-C-N	-5.61	112.91	122.11
1	H	106	THR	C-N-CA	-5.61	112.91	122.11
1	H	114	PHE	CA-CB-CG	5.40	119.20	113.80
1	F	100	ASP	CA-CB-CG	5.21	117.81	112.60

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1018	0	978	36	0
1	B	1088	0	1049	29	0
1	C	1032	0	1002	41	0
1	D	1046	0	1018	35	0
1	E	1029	0	1002	43	0
1	F	1033	0	1004	38	0
1	G	1025	0	998	30	0
1	H	1016	0	981	39	0
2	A	19	18	0	1	0
2	C	19	18	0	1	0
2	E	19	18	0	0	0
2	F	19	18	0	0	0
2	G	19	18	0	0	0
2	H	19	18	0	0	0
3	B	7	10	10	0	0
4	F	6	8	8	0	0
4	H	12	16	16	1	0
5	A	36	0	0	1	0
5	B	40	0	0	0	0
5	C	38	0	0	2	0
5	D	41	0	0	0	0
5	E	32	0	0	1	0
5	F	35	0	0	0	0
5	G	23	0	0	1	0
5	H	44	0	0	2	0
All	All	8715	142	8066	240	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 15.

All (240) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:104:LEU:HD23	1:A:108:ILE:HD12	1.31	1.12
1:A:39:MET:HG2	1:A:107:THR:HG22	1.24	1.09

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:39:MET:HG2	1:C:107:THR:CG2	1.95	0.96
1:C:39:MET:HG2	1:C:107:THR:HG22	1.46	0.94
1:F:13:TRP:HH2	1:G:69:MET:HE3	1.33	0.90
1:A:39:MET:HG2	1:A:107:THR:CG2	2.01	0.90
1:E:69:MET:HE3	1:E:77:VAL:HG22	1.52	0.89
1:A:15:ILE:HD12	1:A:40:PBF:HK1	1.54	0.87
1:F:104:LEU:HA	1:F:108:ILE:HD12	1.61	0.81
1:B:40:PBF:CK1	1:B:86:ILE:HG23	2.12	0.79
1:E:111:HIS:HA	1:E:114:PHE:CE2	2.19	0.77
1:A:104:LEU:CD2	1:A:108:ILE:HD12	2.14	0.75
1:C:100:ASP:O	1:C:104:LEU:HD23	1.87	0.75
1:D:12:CYS:SG	1:D:85:GLU:HG2	2.28	0.74
1:A:15:ILE:HD12	1:A:40:PBF:CK1	2.18	0.73
1:C:98:GLN:O	1:C:104:LEU:HD21	1.87	0.73
1:C:13:TRP:HH2	1:D:69:MET:HE2	1.53	0.73
1:F:119:ARG:HG2	1:F:119:ARG:HH11	1.55	0.70
1:D:51:ARG:NH2	1:D:129:GLU:OE2	2.22	0.70
1:B:103:PHE:CE2	1:B:108:ILE:HD11	2.26	0.70
1:F:13:TRP:HD1	1:F:86:ILE:HD11	1.57	0.69
1:C:13:TRP:HH2	1:D:69:MET:CE	2.05	0.69
1:G:58:ASP:HB2	5:G:308:HOH:O	1.93	0.69
1:D:40:PBF:CK2	1:D:86:ILE:HG23	2.23	0.68
1:C:13:TRP:CH2	1:D:69:MET:CE	2.77	0.68
1:A:85:GLU:OE2	1:B:55:LYS:NZ	2.27	0.68
1:B:91:ILE:HD12	1:B:95:GLN:CD	2.20	0.67
1:F:30:ASN:OD1	1:F:34:LYS:HD2	1.95	0.67
1:H:16:TYR:HB3	1:H:80:TYR:CD2	2.30	0.67
1:F:121:LYS:HE3	1:G:75:GLU:O	1.95	0.66
1:F:13:TRP:CH2	1:G:69:MET:HE3	2.24	0.66
1:B:15:ILE:HG12	1:B:40:PBF:CK2	2.26	0.66
1:A:90:ASP:HB2	1:A:92:GLU:OE1	1.97	0.65
1:F:69:MET:HE3	1:G:13:TRP:HH2	1.61	0.65
1:G:35:VAL:O	1:G:38:PRO:HD2	1.97	0.65
1:E:90:ASP:HB2	1:E:92:GLU:OE1	1.97	0.65
1:D:31:HIS:CE1	1:D:35:VAL:HG11	2.32	0.65
1:E:69:MET:CE	1:E:77:VAL:HG22	2.27	0.64
1:E:112:PHE:CZ	1:F:105:ARG:HB2	2.33	0.64
1:E:46:LYS:HD3	1:E:47:TYR:CE2	2.33	0.64
1:H:47:TYR:CE1	1:H:96:THR:HG22	2.32	0.64
1:G:75:GLU:H	1:G:75:GLU:CD	2.06	0.63
1:A:39:MET:CG	1:A:107:THR:CG2	2.74	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:73:PRO:HB2	1:E:75:GLU:OE1	1.99	0.62
1:E:92:GLU:O	1:E:96:THR:HG23	1.98	0.62
1:F:19:LYS:HE2	1:F:23:GLN:O	2.00	0.62
1:D:51:ARG:HH22	1:D:129:GLU:CD	2.08	0.62
1:F:119:ARG:HG2	1:F:119:ARG:NH1	2.13	0.61
1:A:15:ILE:CD1	1:A:40:PBF:HK1	2.29	0.60
1:C:39:MET:HG2	1:C:107:THR:HG23	1.84	0.60
1:G:40:PBF:CK1	1:G:86:ILE:HG23	2.32	0.60
1:C:41:ILE:HG23	1:D:132:PHE:HE1	1.67	0.59
1:B:13:TRP:CE3	1:B:123:SER:O	2.56	0.59
1:C:41:ILE:HG23	1:D:132:PHE:CE1	2.38	0.58
1:G:31:HIS:CE1	1:G:35:VAL:HG11	2.37	0.58
1:H:111:HIS:HA	1:H:114:PHE:CE2	2.39	0.58
1:G:37:ALA:O	1:G:41:ILE:HD12	2.03	0.58
1:H:20:LYS:HD2	1:H:112:PHE:O	2.04	0.58
1:C:53:THR:OG1	1:D:129:GLU:HG3	2.04	0.57
1:B:15:ILE:HG12	1:B:40:PBF:HK2	1.86	0.57
1:F:111:HIS:HA	1:F:114:PHE:CE2	2.40	0.57
1:B:37:ALA:HB3	1:B:38:PRO:HD3	1.86	0.57
1:C:40:PBF:HI1	1:C:40:PBF:HE1	1.87	0.57
1:D:103:PHE:CE2	1:D:108:ILE:HD11	2.40	0.57
1:E:127:ILE:HG22	1:H:55:LYS:HG3	1.87	0.56
1:C:13:TRP:CZ3	1:D:69:MET:HE3	2.40	0.56
1:B:13:TRP:HE3	1:B:123:SER:O	1.87	0.56
1:F:13:TRP:HB3	1:F:86:ILE:HG13	1.87	0.56
1:E:37:ALA:HB3	1:E:38:PRO:HD3	1.87	0.56
1:C:108:ILE:O	1:C:111:HIS:HD2	1.89	0.56
1:D:37:ALA:HB3	1:D:38:PRO:CD	2.35	0.55
1:H:9:ARG:N	5:H:302:HOH:O	2.39	0.55
1:F:13:TRP:HH2	1:G:69:MET:CE	2.15	0.55
1:C:28:HIS:O	1:C:32:VAL:HG23	2.05	0.55
1:B:91:ILE:HD12	1:B:95:GLN:OE1	2.05	0.55
1:C:40:PBF:HI1	1:C:40:PBF:CE1	2.37	0.55
1:C:111:HIS:ND1	5:C:301:HOH:O	2.32	0.55
1:D:15:ILE:HG13	1:D:40:PBF:HK1	1.89	0.54
1:C:40:PBF:HE1	1:C:40:PBF:CI1	2.38	0.54
1:A:40:PBF:HE2	1:A:40:PBF:CI2	2.38	0.54
1:F:40:PBF:HE1	1:F:40:PBF:HI2	1.89	0.54
1:C:16:TYR:CE1	1:D:78:LEU:HD22	2.42	0.54
1:D:69:MET:SD	1:D:77:VAL:HG22	2.48	0.54
1:A:108:ILE:HA	2:A:201:A1EAB:C3	2.37	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:20:LYS:HD3	1:H:113:ASN:C	2.33	0.53
1:A:108:ILE:HB	1:A:109:PRO:HD3	1.90	0.53
1:G:73:PRO:HB2	1:G:75:GLU:OE1	2.09	0.53
1:B:15:ILE:HG22	1:B:17:VAL:HG23	1.90	0.53
1:H:39:MET:HE2	1:H:107:THR:HG22	1.91	0.53
1:C:108:ILE:HA	2:C:201:A1EAB:C3	2.39	0.53
1:C:127:ILE:HG22	1:D:55:LYS:HG3	1.91	0.53
1:F:40:PBF:HE1	1:F:40:PBF:CI2	2.39	0.52
1:E:124:LEU:HB3	1:H:65:GLN:HB2	1.92	0.52
1:B:40:PBF:CE1	1:B:40:PBF:HI1	2.39	0.52
1:E:127:ILE:HG22	1:H:55:LYS:CG	2.40	0.52
1:E:128:GLU:OE2	1:H:33:SER:OG	2.23	0.52
1:G:40:PBF:HK1	1:G:86:ILE:HG23	1.91	0.52
1:G:30:ASN:OD1	1:G:34:LYS:HD2	2.10	0.52
1:D:55:LYS:HB2	1:D:83:VAL:HB	1.92	0.52
1:F:127:ILE:O	1:F:127:ILE:HG13	2.10	0.51
1:C:111:HIS:CE1	5:C:301:HOH:O	2.63	0.51
1:D:40:PBF:HK2	1:D:86:ILE:HG23	1.91	0.51
1:F:47:TYR:HE1	1:F:96:THR:HG22	1.73	0.51
1:C:39:MET:O	1:C:107:THR:HG21	2.10	0.51
1:A:104:LEU:HD23	1:A:108:ILE:CD1	2.22	0.51
1:F:40:PBF:HI2	1:F:40:PBF:CE1	2.41	0.51
1:H:106:THR:O	4:H:202:GOL:H32	2.11	0.51
1:C:124:LEU:C	1:C:124:LEU:HD13	2.35	0.51
1:A:16:TYR:HB3	1:A:80:TYR:CD2	2.45	0.50
1:A:124:LEU:HB3	1:B:65:GLN:HB2	1.91	0.50
1:A:127:ILE:HG22	1:B:55:LYS:HG3	1.93	0.50
1:E:91:ILE:HG21	1:H:64:LYS:HB3	1.93	0.50
1:E:96:THR:HG22	1:G:71:GLY:HA2	1.92	0.50
1:D:20:LYS:HD2	1:D:113:ASN:C	2.36	0.50
1:A:111:HIS:HA	1:A:114:PHE:CE1	2.46	0.50
1:H:47:TYR:HE1	1:H:96:THR:HG22	1.77	0.50
1:E:75:GLU:CD	1:E:75:GLU:H	2.20	0.49
1:E:104:LEU:HD21	1:F:108:ILE:CG2	2.42	0.49
1:A:39:MET:CG	1:A:107:THR:HG22	2.15	0.49
1:A:58:ASP:OD1	1:A:58:ASP:C	2.53	0.49
1:E:40:PBF:HE1	1:E:40:PBF:CI1	2.42	0.49
1:C:13:TRP:CH2	1:D:69:MET:HE2	2.40	0.49
1:E:78:LEU:HD22	1:H:16:TYR:CE1	2.47	0.49
1:A:51:ARG:O	1:A:86:ILE:HA	2.13	0.48
1:H:51:ARG:O	1:H:86:ILE:HA	2.12	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:122:GLY:HA3	1:D:69:MET:HE1	1.96	0.48
1:E:36:ASN:O	1:E:37:ALA:C	2.57	0.48
1:E:51:ARG:O	1:E:86:ILE:HA	2.13	0.48
1:E:75:GLU:O	1:H:121:LYS:HE3	2.12	0.48
1:H:23:GLN:NE2	1:H:113:ASN:O	2.43	0.48
1:C:65:GLN:HB2	1:D:124:LEU:HB3	1.95	0.48
1:E:84:PHE:N	1:E:84:PHE:CD1	2.81	0.48
1:B:20:LYS:HD2	1:B:112:PHE:O	2.14	0.47
1:D:62:LYS:N	1:D:63:PRO:CD	2.76	0.47
1:C:37:ALA:HB3	1:C:38:PRO:HD3	1.96	0.47
1:B:58:ASP:C	1:B:58:ASP:OD1	2.58	0.47
1:F:78:LEU:HB2	1:G:123:SER:HB2	1.97	0.47
1:F:49:ILE:O	1:G:132:PHE:HD2	1.97	0.47
1:F:49:ILE:HG22	1:G:132:PHE:HE2	1.79	0.47
1:A:16:TYR:O	1:A:120:SER:HB3	2.15	0.47
1:A:55:LYS:HB2	1:A:83:VAL:HB	1.97	0.47
1:B:13:TRP:CZ2	1:B:15:ILE:HD11	2.51	0.46
1:F:39:MET:HG3	1:F:107:THR:HG22	1.95	0.46
1:G:99:LYS:HD2	1:G:99:LYS:HA	1.68	0.46
1:E:132:PHE:CD1	1:H:45:LYS:HG2	2.50	0.46
1:F:78:LEU:HD23	1:F:78:LEU:HA	1.76	0.46
1:G:55:LYS:HB2	1:G:83:VAL:HB	1.96	0.46
1:H:40:PBF:CI2	1:H:40:PBF:HE1	2.45	0.46
1:E:55:LYS:NZ	1:E:85:GLU:OE2	2.46	0.46
1:F:104:LEU:O	1:F:109:PRO:HD3	2.16	0.46
1:E:9:ARG:HG2	1:E:126:TRP:CE3	2.51	0.46
1:B:104:LEU:O	1:B:109:PRO:HD3	2.16	0.46
1:D:40:PBF:CI2	1:D:86:ILE:HG23	2.46	0.46
1:E:108:ILE:N	1:E:109:PRO:CD	2.79	0.46
1:H:99:LYS:HA	1:H:99:LYS:HD3	1.70	0.46
1:A:40:PBF:HI2	1:A:40:PBF:CE2	2.46	0.45
1:H:89:LYS:HG3	5:H:314:HOH:O	2.15	0.45
1:B:102:GLU:OE1	1:B:105:ARG:NE	2.43	0.45
1:E:14:SER:HA	1:E:84:PHE:O	2.16	0.45
1:E:16:TYR:CE1	1:H:78:LEU:HD22	2.51	0.45
1:H:124:LEU:C	1:H:124:LEU:HD13	2.42	0.45
1:H:111:HIS:HA	1:H:114:PHE:CZ	2.51	0.45
1:A:40:PBF:CI2	1:A:40:PBF:CE2	2.93	0.45
1:E:16:TYR:CZ	1:H:78:LEU:HD22	2.52	0.45
1:F:10:LEU:HD23	1:F:10:LEU:HA	1.66	0.45
1:C:91:ILE:HG21	1:D:64:LYS:HB3	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:109:PRO:O	1:E:112:PHE:HB2	2.17	0.45
1:E:69:MET:HE1	1:H:122:GLY:CA	2.47	0.45
1:A:49:ILE:CD1	1:A:49:ILE:N	2.80	0.45
1:E:58:ASP:OD1	1:E:58:ASP:C	2.60	0.45
1:B:14:SER:HA	1:B:84:PHE:O	2.17	0.44
1:D:58:ASP:C	1:D:58:ASP:OD1	2.58	0.44
1:C:13:TRP:HZ3	1:D:69:MET:HE3	1.81	0.44
1:G:13:TRP:CZ2	1:G:15:ILE:HD11	2.53	0.44
1:C:108:ILE:O	1:C:111:HIS:CD2	2.68	0.44
1:A:117:MET:HA	1:A:117:MET:HE3	1.99	0.44
1:H:31:HIS:CE1	1:H:35:VAL:HG11	2.53	0.44
1:E:94:ILE:O	1:E:98:GLN:HG3	2.17	0.44
1:E:69:MET:HE1	1:H:122:GLY:HA3	2.00	0.43
1:H:14:SER:C	1:H:15:ILE:HD13	2.43	0.43
1:A:108:ILE:O	1:A:111:HIS:CD2	2.72	0.43
1:B:111:HIS:C	1:B:111:HIS:CD2	2.96	0.43
1:E:89:LYS:HD2	5:E:312:HOH:O	2.18	0.43
1:C:47:TYR:CE1	1:C:96:THR:HG22	2.53	0.43
1:D:111:HIS:HA	1:D:114:PHE:CE1	2.53	0.43
1:F:47:TYR:CE1	1:F:96:THR:HG22	2.53	0.43
1:A:41:ILE:HG23	1:B:132:PHE:CZ	2.54	0.43
1:E:111:HIS:CD2	1:E:111:HIS:C	2.96	0.43
1:G:124:LEU:HD13	1:G:124:LEU:C	2.44	0.43
1:A:33:SER:OG	1:B:128:GLU:OE1	2.31	0.43
1:C:20:LYS:HB3	1:C:20:LYS:HE2	1.48	0.42
1:C:88:VAL:HG11	1:C:94:ILE:HD11	2.00	0.42
1:D:10:LEU:HD23	1:D:10:LEU:HA	1.75	0.42
1:C:111:HIS:HA	1:C:114:PHE:CE1	2.54	0.42
1:A:91:ILE:HG21	1:B:64:LYS:HB3	2.00	0.42
1:F:39:MET:CG	1:F:107:THR:HG22	2.49	0.42
1:B:124:LEU:HD13	1:B:124:LEU:C	2.45	0.42
1:C:106:THR:OG1	1:C:107:THR:N	2.53	0.42
1:F:100:ASP:OD1	1:F:102:GLU:HB2	2.20	0.42
1:F:39:MET:HE3	1:F:39:MET:HB2	1.86	0.41
1:H:16:TYR:HB3	1:H:80:TYR:CG	2.56	0.41
1:C:16:TYR:CZ	1:D:78:LEU:HD22	2.55	0.41
1:C:106:THR:O	1:C:109:PRO:HD2	2.21	0.41
1:C:104:LEU:O	1:C:109:PRO:HD3	2.20	0.41
1:D:39:MET:HB2	1:D:39:MET:HE2	1.60	0.41
1:F:124:LEU:HB3	1:G:65:GLN:HB2	2.03	0.41
1:E:51:ARG:CD	1:H:131:THR:HG22	2.51	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:20:LYS:HB3	1:H:20:LYS:HE2	1.75	0.41
1:H:58:ASP:OD1	1:H:58:ASP:C	2.60	0.41
1:C:80:TYR:OH	1:C:121:LYS:HD2	2.20	0.41
1:E:85:GLU:OE1	1:H:55:LYS:NZ	2.28	0.41
1:G:75:GLU:CD	1:G:75:GLU:N	2.76	0.41
1:A:87:ILE:HD11	5:A:329:HOH:O	2.21	0.41
1:H:20:LYS:HA	1:H:21:PRO:HD3	1.93	0.41
1:B:44:LEU:HD13	1:B:52:TYR:HB2	2.02	0.41
1:B:46:LYS:HD3	1:B:47:TYR:CE2	2.55	0.41
1:C:14:SER:HA	1:C:84:PHE:O	2.21	0.41
1:F:16:TYR:CE1	1:G:78:LEU:HD22	2.56	0.41
1:H:124:LEU:HD13	1:H:125:THR:N	2.35	0.41
1:A:37:ALA:HB3	1:A:38:PRO:HD3	2.03	0.41
1:A:43:PHE:O	1:A:47:TYR:HD2	2.04	0.41
1:F:49:ILE:HG13	1:F:88:VAL:HG12	2.03	0.41
1:G:13:TRP:HB3	1:G:86:ILE:HG12	2.03	0.41
1:D:47:TYR:CG	1:D:97:MET:HB2	2.56	0.40
1:E:20:LYS:O	1:E:23:GLN:HB3	2.21	0.40
1:E:41:ILE:N	1:E:42:PRO:CD	2.84	0.40
1:E:45:LYS:O	1:E:46:LYS:C	2.61	0.40
1:G:117:MET:HE3	1:G:117:MET:HB3	1.83	0.40
1:H:55:LYS:O	1:H:82:THR:HA	2.21	0.40
1:B:40:PBF:CE1	1:B:40:PBF:CI1	2.98	0.40
1:D:10:LEU:HD21	1:D:89:LYS:HG3	2.03	0.40
1:A:69:MET:HE2	1:A:69:MET:HB3	1.92	0.40
1:E:40:PBF:HI1	1:E:40:PBF:CE1	2.52	0.40
1:F:13:TRP:HB3	1:F:86:ILE:CG1	2.51	0.40
1:G:20:LYS:HE3	1:G:20:LYS:HB3	1.78	0.40
1:A:12:CYS:SG	1:B:55:LYS:HE3	2.62	0.40
1:F:49:ILE:HG22	1:G:132:PHE:CE2	2.56	0.40
1:F:106:THR:OG1	1:F:107:THR:HG23	2.21	0.40
1:F:69:MET:HE3	1:G:13:TRP:CH2	2.50	0.40
1:H:40:PBF:HI2	1:H:40:PBF:CE1	2.51	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	121/141 (86%)	118 (98%)	3 (2%)	0	100	100
1	B	130/141 (92%)	127 (98%)	3 (2%)	0	100	100
1	C	122/141 (86%)	118 (97%)	3 (2%)	1 (1%)	16	16
1	D	124/141 (88%)	121 (98%)	3 (2%)	0	100	100
1	E	121/141 (86%)	119 (98%)	1 (1%)	1 (1%)	16	16
1	F	122/141 (86%)	120 (98%)	1 (1%)	1 (1%)	16	16
1	G	121/141 (86%)	118 (98%)	3 (2%)	0	100	100
1	H	121/141 (86%)	118 (98%)	3 (2%)	0	100	100
All	All	982/1128 (87%)	959 (98%)	20 (2%)	3 (0%)	37	42

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	114	PHE
1	C	114	PHE
1	F	114	PHE

5.3.2 Protein sidechains

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	110/127 (87%)	110 (100%)	0	100	100
1	B	118/127 (93%)	116 (98%)	2 (2%)	56	71

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	112/127 (88%)	109 (97%)	3 (3%)	40	53
1	D	113/127 (89%)	112 (99%)	1 (1%)	75	86
1	E	112/127 (88%)	111 (99%)	1 (1%)	75	86
1	F	112/127 (88%)	109 (97%)	3 (3%)	40	53
1	G	111/127 (87%)	110 (99%)	1 (1%)	75	86
1	H	109/127 (86%)	107 (98%)	2 (2%)	54	69
All	All	897/1016 (88%)	884 (99%)	13 (1%)	62	77

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	107	THR
1	B	118	THR
1	C	20	LYS
1	C	34	LYS
1	C	49	ILE
1	D	55	LYS
1	E	102	GLU
1	F	69	MET
1	F	75	GLU
1	F	99	LYS
1	G	90	ASP
1	H	49	ILE
1	H	107	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	98	GLN
1	A	111	HIS
1	B	57	ASN
1	F	72	GLN
1	F	95	GLN
1	G	72	GLN
1	G	76	ASN
1	G	95	GLN
1	H	57	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

8 non-standard protein/DNA/RNA residues 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	$\# Z > 2$
1	PBF	B	40	1	19,20,21	0.40	0	23,26,28	0.39	0
1	PBF	C	40	1	19,20,21	0.27	0	23,26,28	0.83	0
1	PBF	D	40	1	19,20,21	0.43	0	23,26,28	0.52	0
1	PBF	E	40	1	19,20,21	0.23	0	23,26,28	0.73	1 (4%)
1	PBF	G	40	1	19,20,21	0.41	0	23,26,28	0.85	1 (4%)
1	PBF	A	40	1	19,20,21	0.36	0	23,26,28	0.45	0
1	PBF	F	40	1	19,20,21	0.40	0	23,26,28	0.82	0
1	PBF	H	40	1	19,20,21	0.30	0	23,26,28	0.71	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
1	PBF	B	40	1	-	4/13/14/16	0/2/2/2
1	PBF	C	40	1	-	2/13/14/16	0/2/2/2
1	PBF	D	40	1	-	2/13/14/16	0/2/2/2
1	PBF	E	40	1	-	2/13/14/16	0/2/2/2
1	PBF	G	40	1	-	4/13/14/16	0/2/2/2
1	PBF	A	40	1	-	2/13/14/16	0/2/2/2
1	PBF	F	40	1	-	2/13/14/16	0/2/2/2
1	PBF	H	40	1	-	2/13/14/16	0/2/2/2

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	40	PBF	CG-CB-CA	2.54	119.25	114.10
1	G	40	PBF	CB-CA-C	-2.00	107.71	111.47

There are no chirality outliers.

All (20) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	B	40	PBF	C-CA-CB-CG
1	G	40	PBF	CA-CB-CG-CD2
1	G	40	PBF	CA-CB-CG-CD1
1	C	40	PBF	CA-CB-CG-CD2
1	D	40	PBF	CA-CB-CG-CD2
1	D	40	PBF	CA-CB-CG-CD1
1	C	40	PBF	CA-CB-CG-CD1
1	H	40	PBF	CA-CB-CG-CD2
1	H	40	PBF	CA-CB-CG-CD1
1	A	40	PBF	CA-CB-CG-CD2
1	B	40	PBF	CA-CB-CG-CD2
1	A	40	PBF	CA-CB-CG-CD1
1	F	40	PBF	CA-CB-CG-CD2
1	B	40	PBF	CA-CB-CG-CD1
1	F	40	PBF	CA-CB-CG-CD1
1	E	40	PBF	CA-CB-CG-CD1
1	E	40	PBF	CA-CB-CG-CD2
1	B	40	PBF	N-CA-CB-CG
1	G	40	PBF	C-CA-CB-CG
1	G	40	PBF	N-CA-CB-CG

There are no ring outliers.

8 monomers are involved in 27 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	B	40	PBF	5	0
1	C	40	PBF	3	0
1	D	40	PBF	4	0
1	E	40	PBF	2	0
1	G	40	PBF	2	0
1	A	40	PBF	6	0
1	F	40	PBF	3	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	H	40	PBF	2	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

10 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
4	GOL	H	202	-	5,5,5	1.22	1 (20%)	5,5,5	1.93	2 (40%)
2	A1EAB	G	201	-	18,19,19	2.02	1 (5%)	21,26,26	0.41	0
3	PEG	B	201	-	6,6,6	0.41	0	5,5,5	0.17	0
4	GOL	F	201	-	5,5,5	1.03	0	5,5,5	0.85	0
4	GOL	H	201	-	5,5,5	1.22	0	5,5,5	0.51	0
2	A1EAB	H	203	-	18,19,19	1.63	1 (5%)	21,26,26	0.77	1 (4%)
2	A1EAB	A	201	-	18,19,19	1.82	1 (5%)	21,26,26	0.71	0
2	A1EAB	E	201	-	18,19,19	1.98	1 (5%)	21,26,26	0.43	0
2	A1EAB	F	202	-	18,19,19	2.10	1 (5%)	21,26,26	0.43	0
2	A1EAB	C	201	-	18,19,19	2.01	1 (5%)	21,26,26	0.34	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	GOL	H	202	-	-	2/4/4/4	-
2	A1EAB	G	201	-	-	2/10/16/16	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	PEG	B	201	-	-	3/4/4/4	-
4	GOL	F	201	-	-	0/4/4/4	-
4	GOL	H	201	-	-	0/4/4/4	-
2	A1EAB	H	203	-	-	2/10/16/16	0/1/1/1
2	A1EAB	A	201	-	-	3/10/16/16	0/1/1/1
2	A1EAB	E	201	-	-	2/10/16/16	0/1/1/1
2	A1EAB	F	202	-	-	1/10/16/16	0/1/1/1
2	A1EAB	C	201	-	-	2/10/16/16	0/1/1/1

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	202	A1EAB	C9-C8	-8.81	1.37	1.49
2	G	201	A1EAB	C9-C8	-8.48	1.37	1.49
2	C	201	A1EAB	C9-C8	-8.45	1.37	1.49
2	E	201	A1EAB	C9-C8	-8.30	1.38	1.49
2	A	201	A1EAB	C9-C8	-7.65	1.38	1.49
2	H	203	A1EAB	C9-C8	-6.67	1.40	1.49
4	H	202	GOL	O1-C1	2.45	1.52	1.42

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	H	202	GOL	C3-C2-C1	-3.10	99.65	111.70
4	H	202	GOL	O2-C2-C1	2.85	121.67	109.12
2	H	203	A1EAB	C16-C13-C12	2.13	113.31	109.47

There are no chirality outliers.

All (17) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	201	A1EAB	C4-C7-C8-C10
2	C	201	A1EAB	C4-C7-C8-C9
3	B	201	PEG	O1-C1-C2-O2
4	H	202	GOL	O1-C1-C2-C3
4	H	202	GOL	C1-C2-C3-O3
3	B	201	PEG	O2-C3-C4-O4
2	E	201	A1EAB	C4-C7-C8-C10
2	G	201	A1EAB	C4-C7-C8-C10
2	H	203	A1EAB	C4-C7-C8-C10

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Mol	Chain	Res	Type	Atoms
2	A	201	A1EAB	C4-C7-C8-C9
2	E	201	A1EAB	C4-C7-C8-C9
2	F	202	A1EAB	C4-C7-C8-C9
2	G	201	A1EAB	C4-C7-C8-C9
2	H	203	A1EAB	C4-C7-C8-C9
3	B	201	PEG	C4-C3-O2-C2
2	A	201	A1EAB	C10-C12-C13-C14
2	A	201	A1EAB	C10-C12-C13-C16

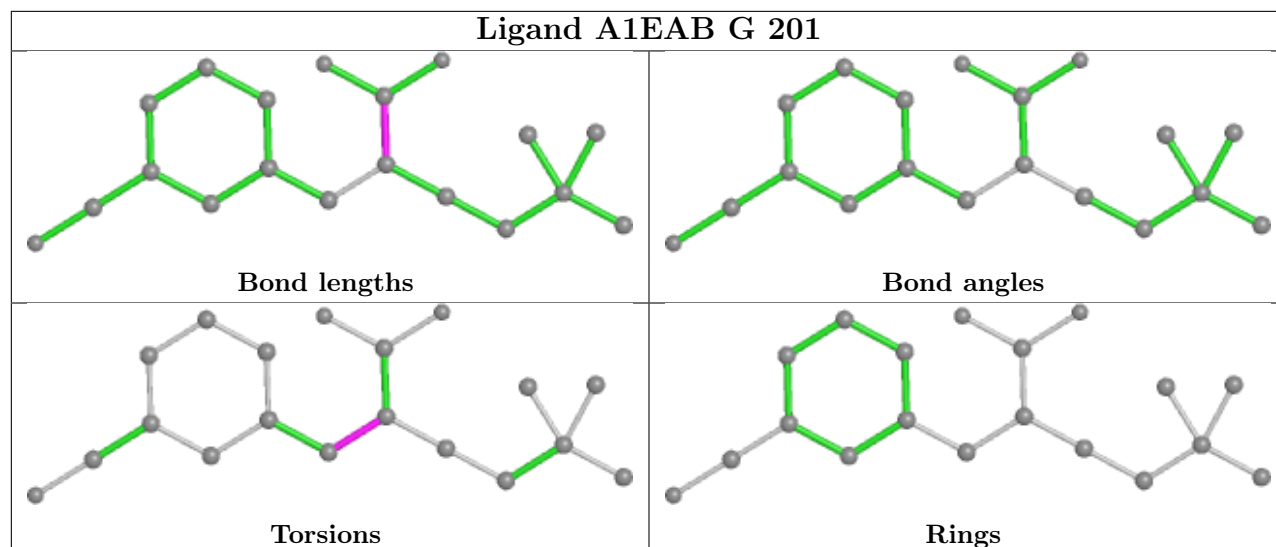
There are no ring outliers.

3 monomers are involved in 3 short contacts:

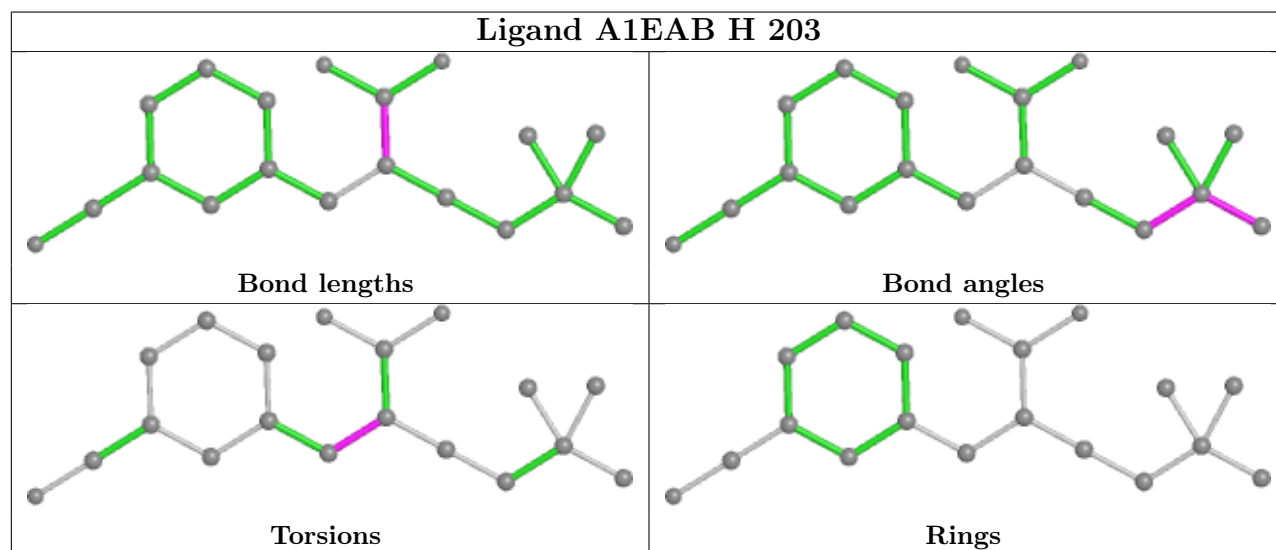
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	H	202	GOL	1	0
2	A	201	A1EAB	1	0
2	C	201	A1EAB	1	0

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 than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

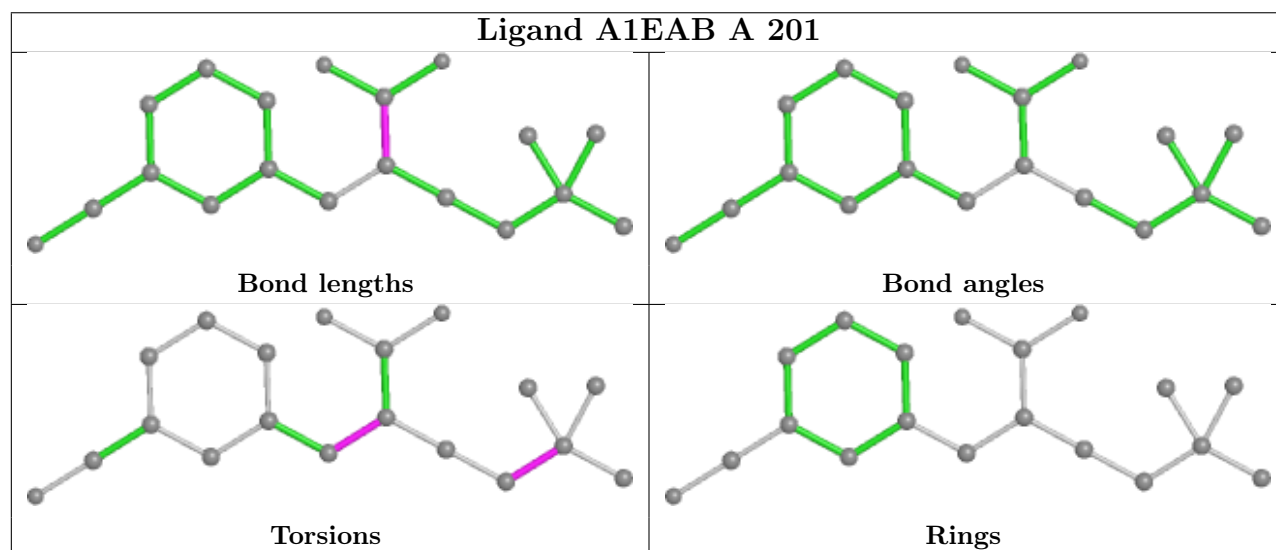
Ligand A1EAB G 201



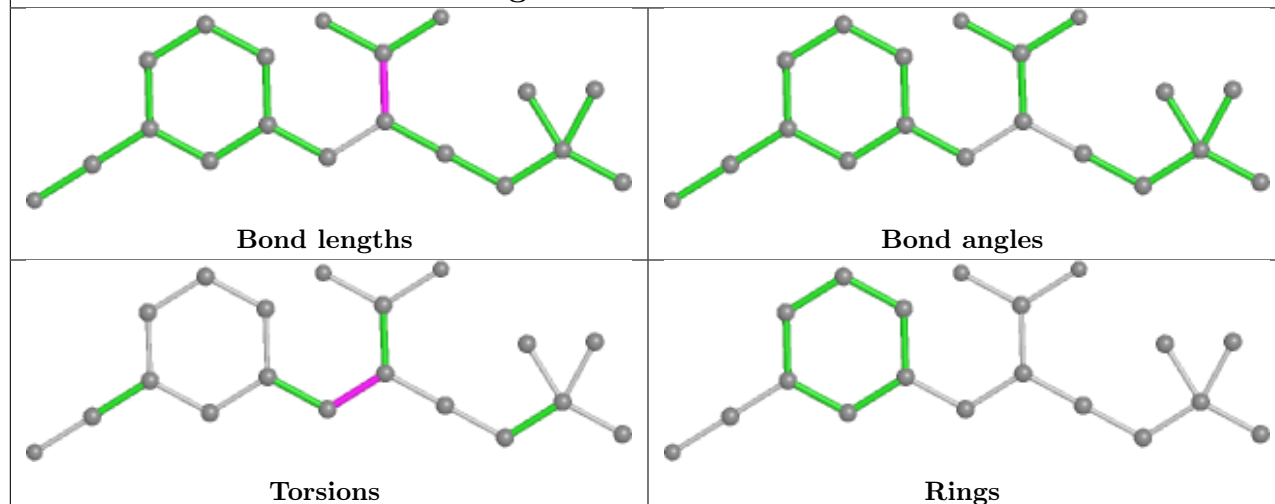
Ligand A1EAB H 203



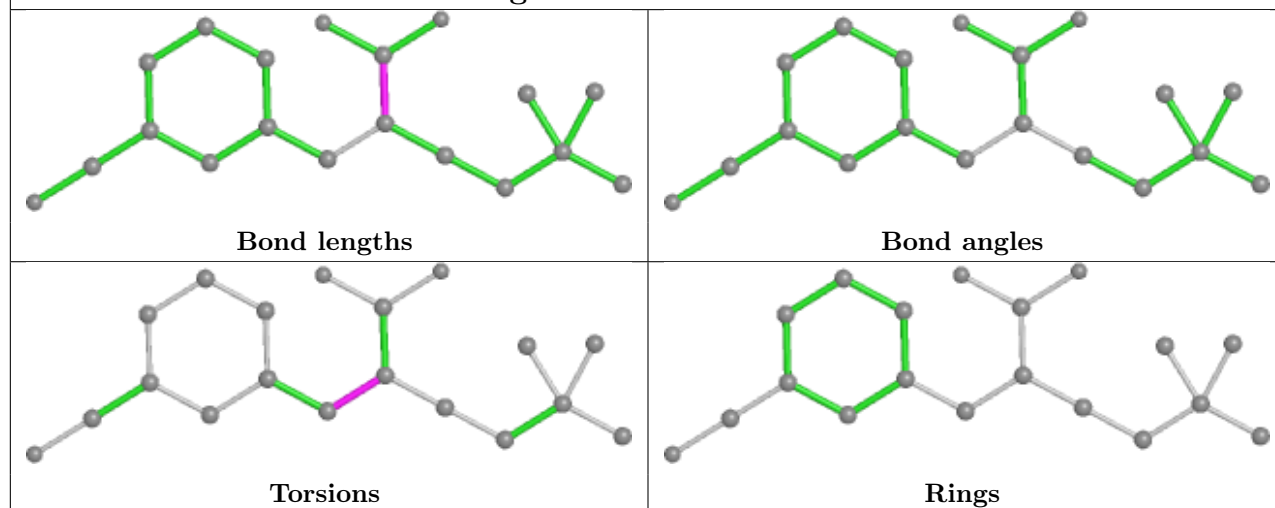
Ligand A1EAB A 201



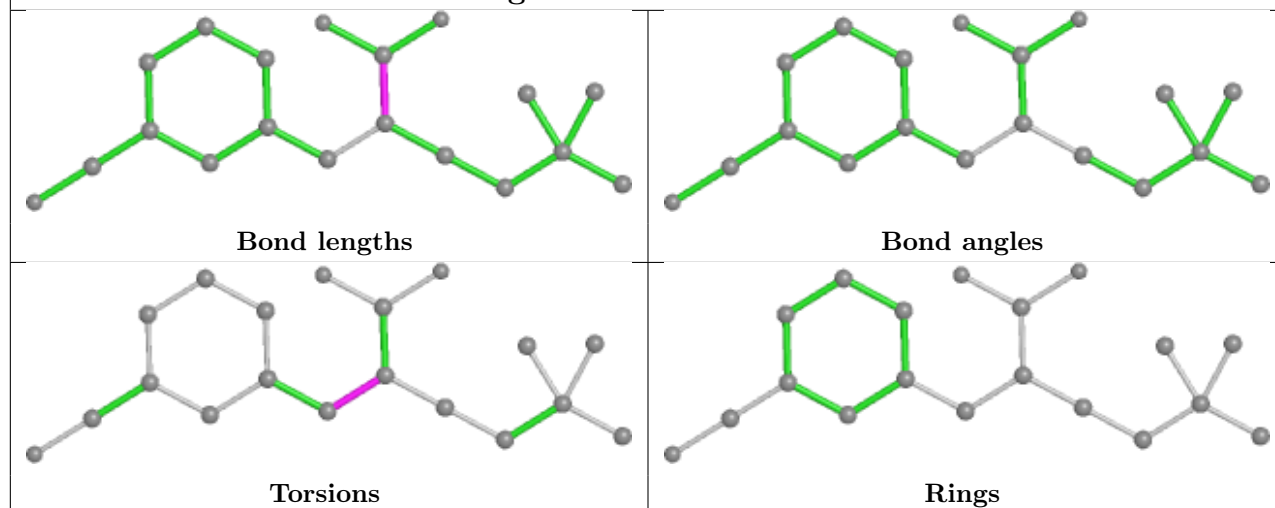
Ligand A1EAB E 201



Ligand A1EAB F 202



Ligand A1EAB C 201



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 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	123/141 (87%)	-0.38	0 100 100	20, 29, 44, 54	0
1	B	132/141 (93%)	-0.29	0 100 100	21, 31, 44, 50	0
1	C	124/141 (87%)	-0.37	0 100 100	22, 30, 42, 56	0
1	D	126/141 (89%)	-0.28	0 100 100	23, 30, 44, 49	0
1	E	123/141 (87%)	-0.21	0 100 100	23, 33, 44, 48	0
1	F	124/141 (87%)	-0.16	0 100 100	21, 30, 42, 54	0
1	G	123/141 (87%)	-0.14	1 (0%) 82 80	23, 34, 44, 54	0
1	H	123/141 (87%)	-0.35	0 100 100	21, 30, 40, 53	0
All	All	998/1128 (88%)	-0.27	1 (0%) 92 91	20, 31, 44, 56	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	22	ASP	2.4

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
1	PBF	B	40	19/20	0.86	0.15	33,35,43,43	0
1	PBF	E	40	19/20	0.86	0.12	27,32,36,36	0
1	PBF	D	40	19/20	0.90	0.14	23,33,40,42	0
1	PBF	A	40	19/20	0.92	0.09	23,29,33,35	0
1	PBF	G	40	19/20	0.93	0.08	25,31,36,37	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
1	PBF	H	40	19/20	0.94	0.07	21,25,30,31	0
1	PBF	C	40	19/20	0.95	0.07	22,28,31,31	0
1	PBF	F	40	19/20	0.95	0.07	23,27,30,32	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

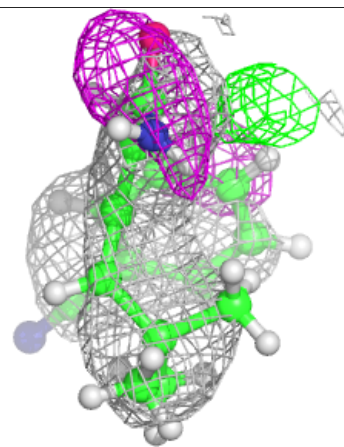
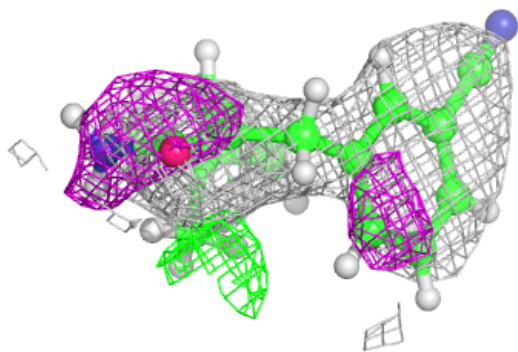
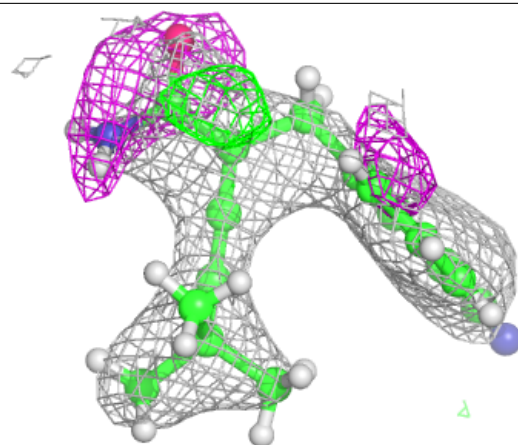
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	A1EAB	C	201	19/19	0.80	0.15	33,45,54,60	0
2	A1EAB	A	201	19/19	0.83	0.12	36,46,56,61	0
2	A1EAB	E	201	19/19	0.85	0.12	34,42,51,51	0
3	PEG	B	201	7/7	0.86	0.15	36,46,56,56	0
2	A1EAB	F	202	19/19	0.87	0.11	32,40,49,50	0
4	GOL	H	201	6/6	0.89	0.12	39,47,54,54	0
4	GOL	H	202	6/6	0.89	0.14	34,41,47,50	0
2	A1EAB	G	201	19/19	0.90	0.10	33,40,49,49	0
2	A1EAB	H	203	19/19	0.93	0.07	28,32,40,40	0
4	GOL	F	201	6/6	0.95	0.11	28,33,40,41	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

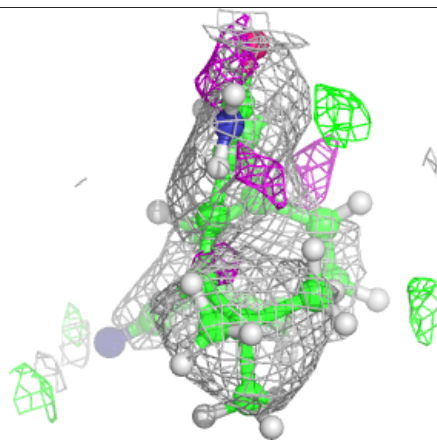
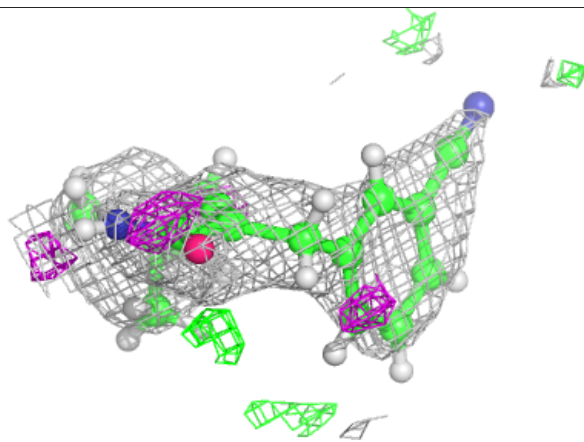
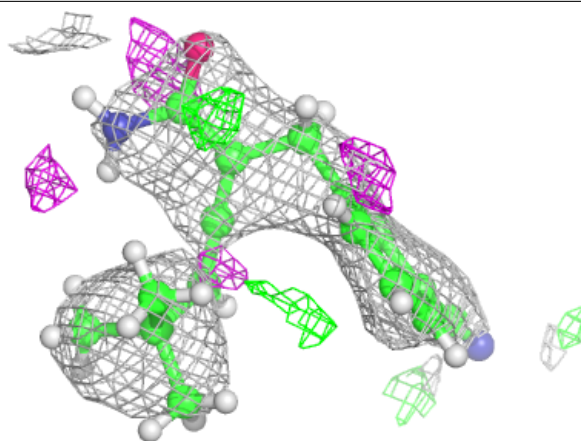
Electron density around A1EAB C 201:

$2mF_o - DF_c$ (at 0.7 rmsd) in gray
 $mF_o - DF_c$ (at 3 rmsd) in purple (negative)
and green (positive)



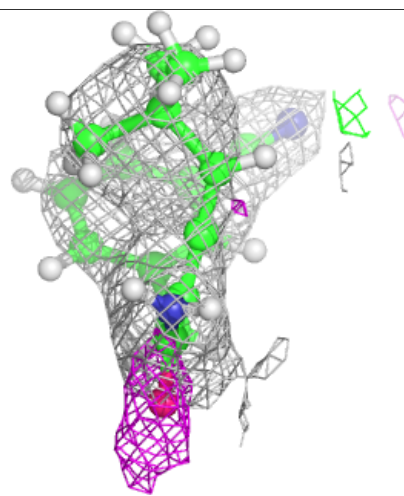
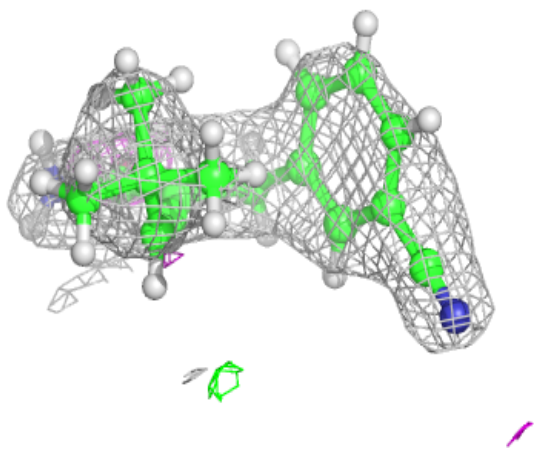
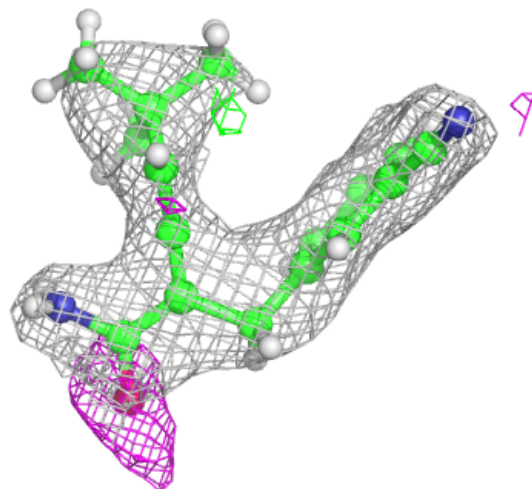
Electron density around A1EAB A 201:

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and green (positive)



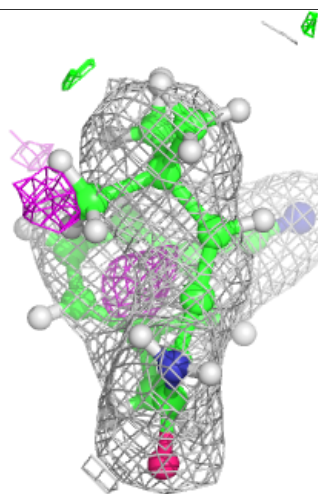
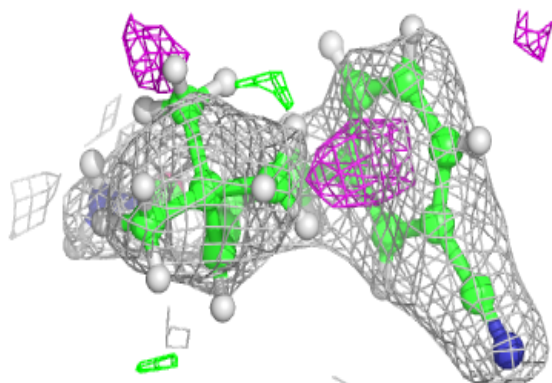
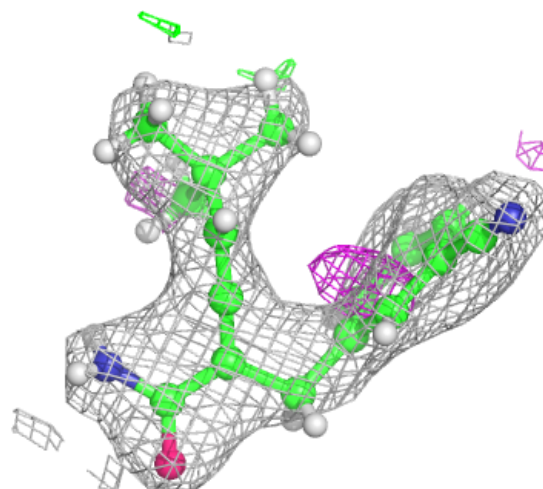
Electron density around A1EAB E 201:

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and green (positive)



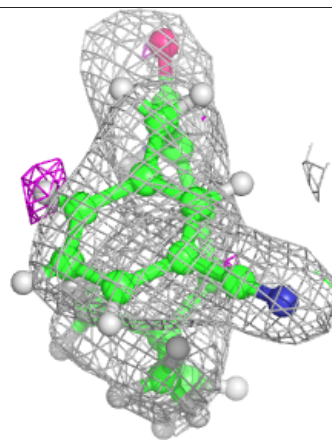
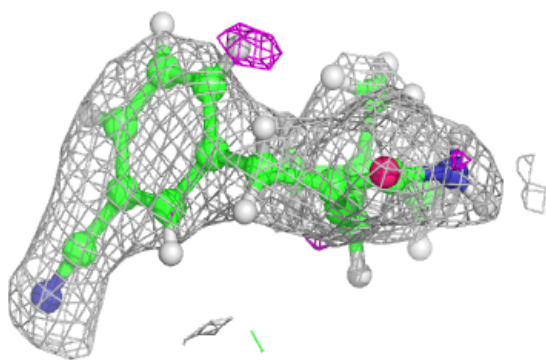
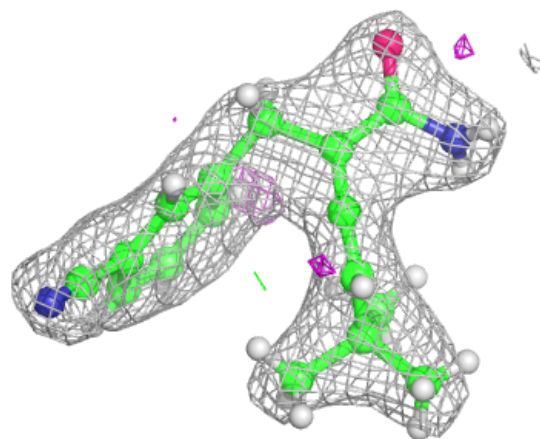
Electron density around A1EAB F 202:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



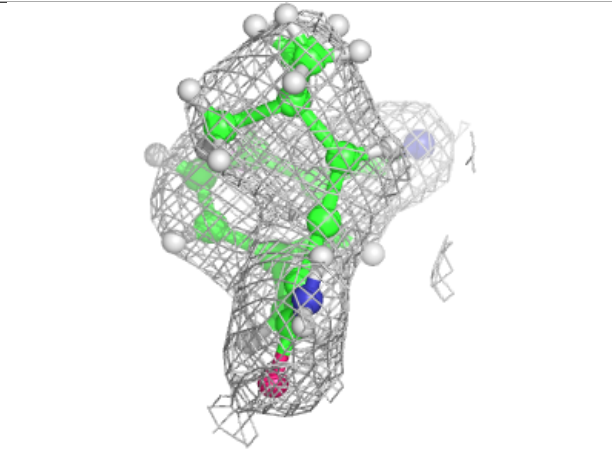
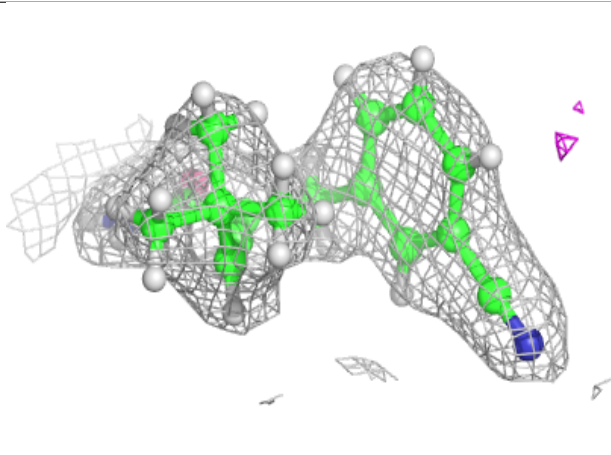
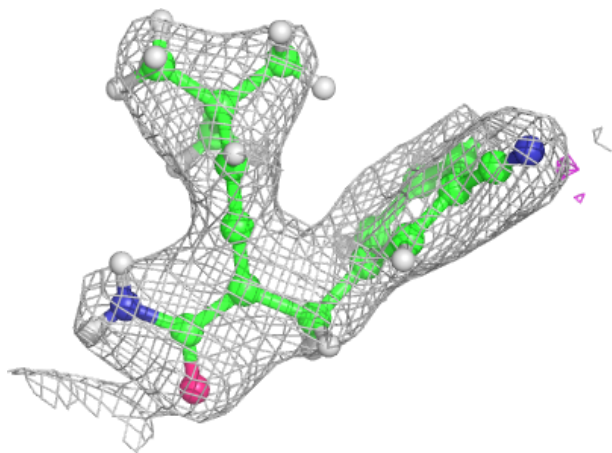
Electron density around A1EAB G 201:

$2mF_o - DF_c$ (at 0.7 rmsd) in gray
 $mF_o - DF_c$ (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around A1EAB H 203:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.