



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 01:44 AM UTC

PDB ID : 6JLG / pdb_00006jlg
Title : Crystal Structure of HasAp with Co-9,10,19,20-Tetraphenylporphycene
Authors : Sakakibara, E.; Shisaka, Y.; Onoda, H.; Sugimoto, H.; Shiro, Y.; Watanabe, Y.; Shoji, O.
Deposited on : 2019-03-05
Resolution : 2.50 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:

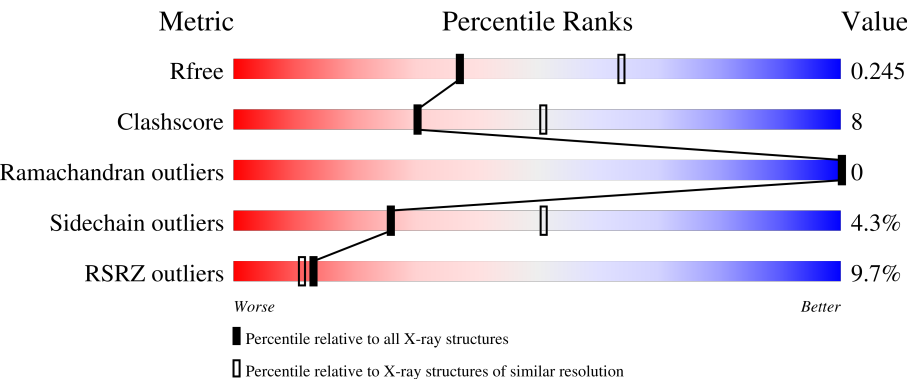
MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

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 2.50 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.



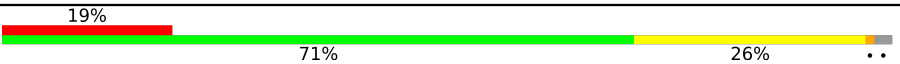
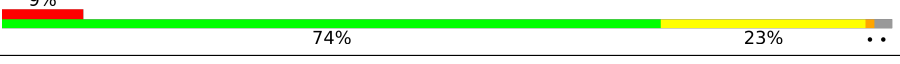
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	184	10% 80% 18% .
1	B	184	9% 80% 17% ..
1	C	184	7% 82% 14% ..
1	D	184	11% 79% 17% ..
1	E	184	5% 80% 18% .

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Mol	Chain	Length	Quality of chain
1	F	184	
1	G	184	
1	H	184	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	BWU	A	201	X	-	-	-
2	BWU	B	201	X	-	-	-
2	BWU	C	201	X	-	-	-
2	BWU	D	201	X	-	-	-
2	BWU	E	201	X	-	-	-
2	BWU	F	201	X	-	-	-
2	BWU	G	201	X	-	-	-
2	BWU	H	201	X	-	-	-

2 Entry composition

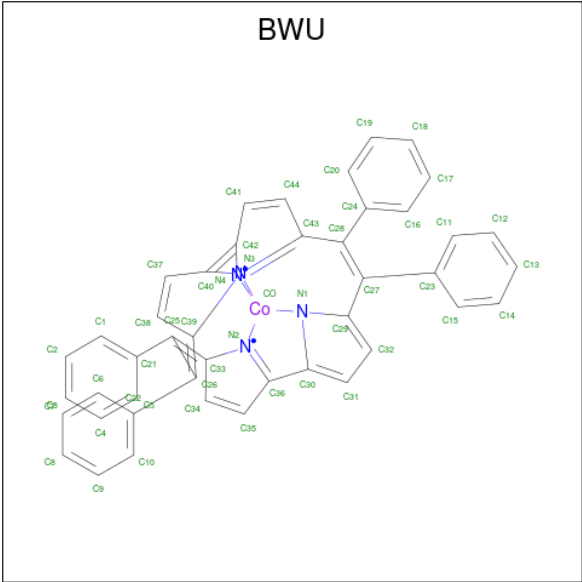
There are 5 unique types of molecules in this entry. The entry contains 11178 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Heme acquisition protein HasA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	181	Total	C	N	O	S	0	0	0
			1316	825	217	273	1			
1	B	181	Total	C	N	O	S	0	0	0
			1316	825	217	273	1			
1	C	177	Total	C	N	O	S	0	0	0
			1290	812	212	265	1			
1	D	181	Total	C	N	O	S	0	0	0
			1316	825	217	273	1			
1	E	181	Total	C	N	O	S	0	0	0
			1316	825	217	273	1			
1	F	181	Total	C	N	O	S	0	0	0
			1316	825	217	273	1			
1	G	181	Total	C	N	O	S	0	0	0
			1316	825	217	273	1			
1	H	181	Total	C	N	O	S	0	0	0
			1316	825	217	273	1			

- Molecule 2 is Tetraphenylporphycene containing cobalt (CCD ID: BWU) (formula: $C_{44}H_{28}CoN_4$).



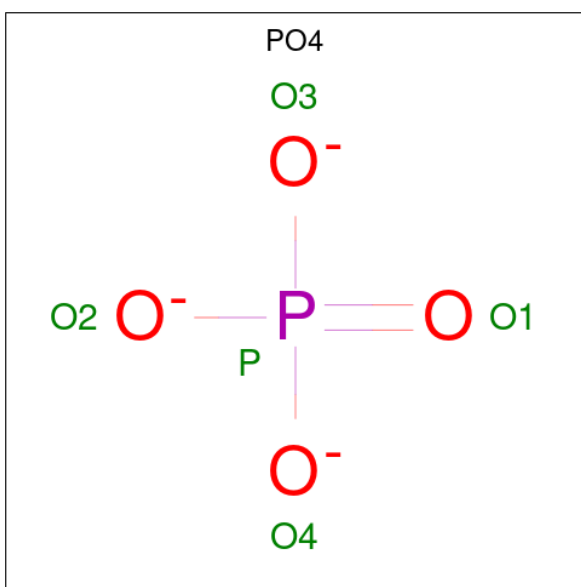
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	Co	N	0	0
			49	44	1	4		
2	B	1	Total	C	Co	N	0	0
			49	44	1	4		
2	C	1	Total	C	Co	N	0	0
			49	44	1	4		
2	D	1	Total	C	Co	N	0	0
			49	44	1	4		
2	E	1	Total	C	Co	N	0	0
			49	44	1	4		
2	F	1	Total	C	Co	N	0	0
			49	44	1	4		
2	G	1	Total	C	Co	N	0	0
			49	44	1	4		
2	H	1	Total	C	Co	N	0	0
			49	44	1	4		

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		
3	D	1	Total	C	O	0	0
			6	3	3		
3	E	1	Total	C	O	0	0
			6	3	3		
3	G	1	Total	C	O	0	0
			6	3	3		
3	G	1	Total	C	O	0	0
			6	3	3		
3	H	1	Total	C	O	0	0
			6	3	3		

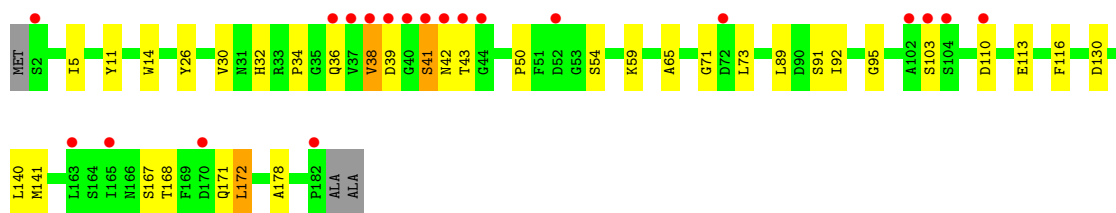
- Molecule 4 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).



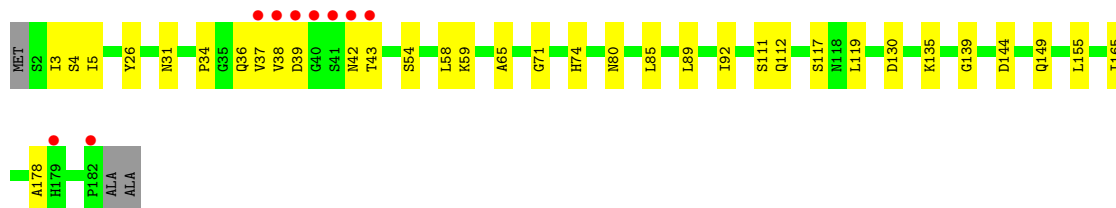
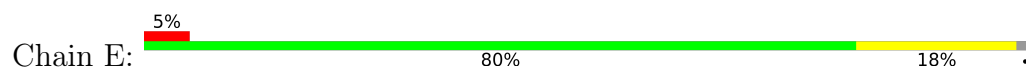
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	E	1	Total	O	P	0	0
			5	4	1		
4	G	1	Total	O	P	0	0
			5	4	1		

- Molecule 5 is water.

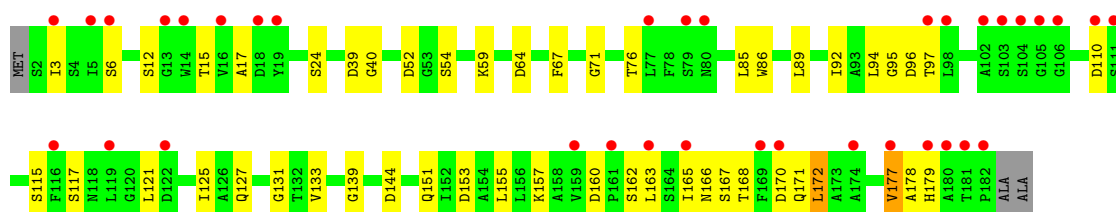
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	32	Total	O	0	0
			32	32		
5	B	36	Total	O	0	0
			36	36		
5	C	18	Total	O	0	0
			18	18		
5	D	17	Total	O	0	0
			17	17		
5	E	47	Total	O	0	0
			47	47		
5	F	11	Total	O	0	0
			11	11		
5	G	41	Total	O	0	0
			41	41		
5	H	18	Total	O	0	0
			18	18		



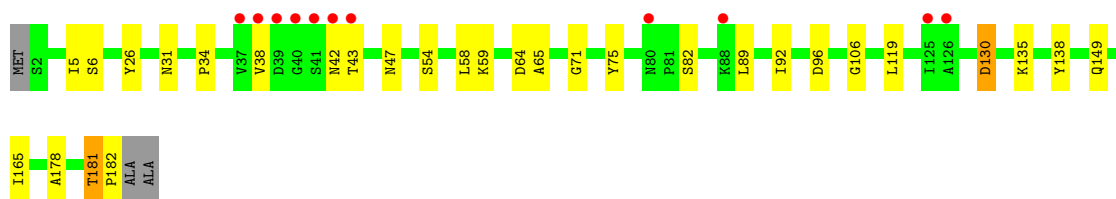
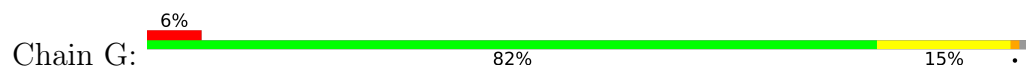
• Molecule 1: Heme acquisition protein HasA



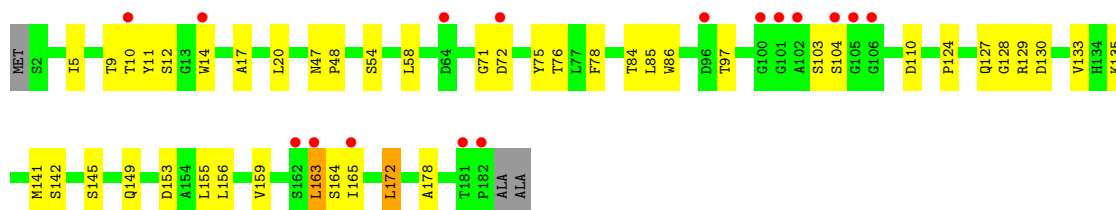
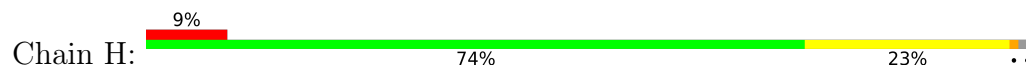
• Molecule 1: Heme acquisition protein HasA



• Molecule 1: Heme acquisition protein HasA



• Molecule 1: Heme acquisition protein HasA



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	237.58Å 92.22Å 93.19Å 90.00° 100.96° 90.00°	Depositor
Resolution (Å)	45.79 – 2.50 45.79 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.7 (45.79-2.50) 99.7 (45.79-2.50)	Depositor EDS
R_{merge}	0.25	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.78 (at 2.51Å)	Xtriage
Refinement program	REFMAC 5.8.0238	Depositor
R, R_{free}	0.213 , 0.241 0.217 , 0.245	Depositor DCC
R_{free} test set	3400 reflections (4.96%)	wwPDB-VP
Wilson B-factor (Å ²)	42.5	Xtriage
Anisotropy	0.436	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 49.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	11178	wwPDB-VP
Average B, all atoms (Å ²)	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 19.00% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: BWU, GOL, PO4

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	0.97	0/1348	1.32	0/1840
1	B	0.97	0/1348	1.31	1/1840 (0.1%)
1	C	0.96	0/1321	1.31	0/1802
1	D	0.98	0/1348	1.31	0/1840
1	E	0.98	0/1348	1.28	0/1840
1	F	0.98	0/1348	1.32	0/1840
1	G	0.98	0/1348	1.27	0/1840
1	H	0.97	0/1348	1.34	0/1840
All	All	0.97	0/10757	1.31	1/14682 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	101	GLY	CA-C-O	-5.22	118.10	122.33

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	A	1316	0	1229	23	0
1	B	1316	0	1229	24	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	1290	0	1209	14	0
1	D	1316	0	1228	26	0
1	E	1316	0	1228	20	0
1	F	1316	0	1229	33	0
1	G	1316	0	1228	17	0
1	H	1316	0	1229	25	0
2	A	49	0	0	0	0
2	B	49	0	0	0	0
2	C	49	0	0	0	0
2	D	49	0	0	1	0
2	E	49	0	0	0	0
2	F	49	0	0	0	0
2	G	49	0	0	1	0
2	H	49	0	0	2	0
3	A	12	0	16	1	0
3	B	12	0	16	0	0
3	D	6	0	8	2	0
3	E	6	0	8	0	0
3	G	12	0	16	1	0
3	H	6	0	8	0	0
4	E	5	0	0	0	0
4	G	5	0	0	1	0
5	A	32	0	0	0	0
5	B	36	0	0	0	0
5	C	18	0	0	0	0
5	D	17	0	0	0	0
5	E	47	0	0	0	0
5	F	11	0	0	0	0
5	G	41	0	0	0	0
5	H	18	0	0	0	0
All	All	11178	0	9881	175	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 8.

All (175) 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:B:3:ILE:HG12	1:B:155:LEU:HB3	1.36	1.04
1:A:112:GLN:NE2	1:A:114:VAL:C	2.33	0.87
1:B:3:ILE:H	1:B:3:ILE:HD12	1.38	0.86

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:42:ASN:O	1:G:42:ASN:ND2	2.09	0.85
1:A:2:SER:OG	3:A:202:GOL:H2	1.77	0.84
1:B:3:ILE:CG1	1:B:155:LEU:HB3	2.07	0.83
1:A:112:GLN:HE22	1:A:114:VAL:C	1.87	0.81
1:C:160:ASP:OD1	1:C:161:PRO:HD2	1.82	0.80
1:H:156:LEU:HB3	1:H:163:LEU:HD12	1.65	0.78
1:F:168:THR:HB	1:F:171:GLN:HG3	1.66	0.77
1:G:43:THR:O	1:G:58:LEU:HD12	1.85	0.77
1:H:129:ARG:O	1:H:135:LYS:HE3	1.90	0.72
1:A:95:GLY:HA3	1:A:109:LEU:HD13	1.73	0.71
1:H:156:LEU:O	1:H:159:VAL:HG12	1.91	0.70
1:D:32:HIS:CE1	1:D:39:ASP:HB3	2.27	0.69
1:F:97:THR:HB	1:F:110:ASP:HB3	1.75	0.68
1:A:97:THR:HB	1:A:110:ASP:HB3	1.77	0.65
1:A:112:GLN:NE2	1:A:114:VAL:O	2.30	0.65
1:F:139:GLY:C	1:F:144:ASP:O	2.40	0.65
1:F:6:SER:HB2	1:F:115:SER:OG	1.97	0.64
1:D:34:PRO:HG2	1:E:34:PRO:HD2	1.79	0.64
1:A:181:THR:HB	1:A:182:PRO:HD2	1.81	0.63
1:F:3:ILE:HD12	1:F:117:SER:O	2.00	0.62
1:D:38:VAL:HG21	1:D:42:ASN:ND2	2.15	0.61
1:E:38:VAL:HG22	1:E:39:ASP:H	1.65	0.61
1:D:36:GLN:HG2	1:D:38:VAL:O	2.00	0.61
1:A:54:SER:HA	1:A:71:GLY:O	2.00	0.60
1:F:153:ASP:OD1	1:F:165:ILE:N	2.35	0.59
1:B:78:PHE:CE2	3:D:202:GOL:H12	2.37	0.59
1:H:153:ASP:HB2	1:H:165:ILE:HD11	1.84	0.59
1:H:20:LEU:HD12	1:H:149:GLN:HG2	1.85	0.58
1:A:181:THR:O	1:A:182:PRO:C	2.45	0.58
1:H:163:LEU:HD11	1:H:172:LEU:HD21	1.86	0.58
1:C:181:THR:HG23	1:C:182:PRO:HD2	1.84	0.57
1:B:78:PHE:HZ	1:D:50:PRO:HB3	1.69	0.56
1:C:48:PRO:HD2	1:C:53:GLY:HA2	1.87	0.56
1:F:15:THR:HG23	1:F:166:ASN:HA	1.87	0.56
1:A:31:ASN:HB3	1:A:36:GLN:OE1	2.05	0.56
1:G:149:GLN:HG3	1:G:165:ILE:HG12	1.87	0.56
1:E:89:LEU:HD21	1:E:92:ILE:HD11	1.88	0.55
1:B:3:ILE:H	1:B:3:ILE:CD1	2.08	0.55
1:E:43:THR:O	1:E:59:LYS:N	2.36	0.55
1:F:172:LEU:O	1:F:177:VAL:HG12	2.06	0.55
1:E:38:VAL:HG22	1:E:39:ASP:N	2.22	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:77:LEU:O	1:B:82:SER:HA	2.06	0.54
1:D:168:THR:O	1:D:172:LEU:HD22	2.08	0.54
1:E:26:TYR:CD2	1:E:65:ALA:HB2	2.43	0.54
1:D:89:LEU:HD21	1:D:92:ILE:HD11	1.90	0.54
1:H:97:THR:HB	1:H:110:ASP:HB3	1.90	0.54
1:A:76:THR:HG21	1:A:125:ILE:CD1	2.37	0.53
1:B:78:PHE:CZ	3:D:202:GOL:H12	2.44	0.53
1:C:116:PHE:HB3	1:C:119:LEU:HD11	1.89	0.53
1:D:95:GLY:HA2	1:D:113:GLU:HG3	1.91	0.53
1:C:36:GLN:O	1:C:36:GLN:HG2	2.08	0.53
1:F:139:GLY:CA	1:F:144:ASP:O	2.57	0.52
1:H:11:TYR:HA	1:H:14:TRP:CD1	2.44	0.51
1:C:160:ASP:OD1	1:C:161:PRO:CD	2.54	0.51
1:E:3:ILE:CG2	1:E:155:LEU:HD13	2.41	0.51
1:E:43:THR:HG23	1:E:58:LEU:HD11	1.92	0.51
1:A:97:THR:O	1:A:98:LEU:C	2.54	0.50
1:B:52:ASP:OD2	1:B:74:HIS:ND1	2.35	0.50
1:C:54:SER:HA	1:C:71:GLY:O	2.11	0.50
1:F:89:LEU:HD21	1:F:92:ILE:HD11	1.92	0.50
1:D:32:HIS:HA	1:D:36:GLN:HE22	1.76	0.50
1:B:153:ASP:OD2	1:B:157:LYS:HE2	2.11	0.50
1:H:54:SER:HA	1:H:71:GLY:O	2.10	0.50
1:H:124:PRO:HG2	1:H:127:GLN:HG3	1.92	0.50
1:D:26:TYR:CD2	1:D:65:ALA:HB2	2.47	0.50
1:F:95:GLY:C	1:F:96:ASP:OD1	2.55	0.50
1:G:59:LYS:HE2	1:G:64:ASP:HA	1.93	0.49
1:B:97:THR:HB	1:B:110:ASP:HB3	1.94	0.49
1:B:129:ARG:HD3	1:G:26:TYR:CG	2.46	0.49
1:H:145:SER:O	1:H:149:GLN:HG3	2.12	0.49
1:B:5:ILE:O	1:B:178:ALA:HA	2.13	0.49
1:E:36:GLN:O	1:E:36:GLN:CG	2.61	0.49
1:H:76:THR:HG23	1:H:86:TRP:HE1	1.78	0.49
1:B:54:SER:HA	1:B:71:GLY:O	2.13	0.49
1:F:76:THR:HG21	1:F:125:ILE:CD1	2.43	0.49
1:A:48:PRO:HD2	1:A:53:GLY:HA2	1.94	0.48
1:F:157:LYS:HA	1:F:160:ASP:O	2.13	0.48
1:C:69:ALA:HB1	1:C:89:LEU:HD11	1.95	0.48
1:E:5:ILE:O	1:E:178:ALA:HA	2.14	0.48
1:F:54:SER:HA	1:F:71:GLY:O	2.14	0.48
1:B:91:SER:HA	1:B:116:PHE:O	2.12	0.48
1:G:130:ASP:O	1:G:135:LYS:NZ	2.46	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:42:ASN:OD1	1:H:78:PHE:CE1	2.66	0.48
1:F:127:GLN:HE21	1:F:131:GLY:HA2	1.77	0.48
1:D:92:ILE:HD13	1:D:140:LEU:HD11	1.95	0.47
1:G:26:TYR:CD2	1:G:65:ALA:HB2	2.49	0.47
1:G:82:SER:OG	4:G:202:PO4:O1	2.28	0.47
1:H:163:LEU:HD13	1:H:164:SER:N	2.30	0.47
1:D:54:SER:HA	1:D:71:GLY:O	2.15	0.47
1:H:17:ALA:N	1:H:165:ILE:HG22	2.30	0.47
1:C:155:LEU:O	1:C:158:ALA:HB3	2.15	0.47
1:A:95:GLY:CA	1:A:109:LEU:HD13	2.43	0.46
1:D:41:SER:HB3	1:D:59:LYS:HB3	1.97	0.46
1:G:5:ILE:O	1:G:178:ALA:HA	2.15	0.46
1:G:89:LEU:HD21	1:G:92:ILE:HD11	1.96	0.46
1:D:168:THR:O	1:D:172:LEU:CD2	2.64	0.46
1:F:76:THR:HG21	1:F:125:ILE:HD12	1.97	0.46
1:F:178:ALA:C	1:F:179:HIS:HD2	2.23	0.46
1:D:30:VAL:HG22	1:D:141:MET:O	2.15	0.46
1:E:149:GLN:CG	1:E:165:ILE:HG12	2.46	0.46
1:F:67:PHE:HB2	1:F:94:LEU:HD23	1.98	0.46
1:E:54:SER:HA	1:E:71:GLY:O	2.16	0.46
1:D:167:SER:HB2	1:D:172:LEU:HD21	1.98	0.46
1:B:3:ILE:HG12	1:B:155:LEU:CB	2.25	0.45
1:F:160:ASP:OD1	1:F:162:SER:N	2.43	0.45
1:A:76:THR:HG21	1:A:125:ILE:HD13	1.98	0.45
1:D:34:PRO:CG	1:E:34:PRO:HD2	2.46	0.45
1:H:9:THR:O	1:H:12:SER:OG	2.28	0.45
1:E:4:SER:OG	1:E:117:SER:HB3	2.17	0.45
1:B:121:LEU:HD11	1:B:151:GLN:HG2	1.98	0.44
1:F:167:SER:HB2	1:F:172:LEU:HD21	1.99	0.44
1:F:15:THR:HG22	1:F:17:ALA:H	1.82	0.44
1:F:76:THR:HG23	1:F:86:TRP:HE1	1.82	0.44
1:B:3:ILE:HD12	1:B:3:ILE:N	2.19	0.44
1:C:5:ILE:O	1:C:178:ALA:HA	2.18	0.44
1:H:72:ASP:N	1:H:72:ASP:OD1	2.51	0.44
1:B:160:ASP:N	1:B:160:ASP:OD1	2.48	0.44
1:F:15:THR:HG22	1:F:17:ALA:N	2.32	0.44
1:A:163:LEU:HD22	1:A:172:LEU:HD23	1.99	0.44
1:C:34:PRO:HG2	1:G:34:PRO:HD2	2.00	0.44
1:D:11:TYR:CE2	1:D:113:GLU:HB3	2.53	0.44
1:E:36:GLN:O	1:E:36:GLN:HG3	2.18	0.44
1:G:138:TYR:HE2	3:G:204:GOL:H31	1.83	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:48:PRO:HD2	1:B:53:GLY:HA2	1.99	0.43
1:C:97:THR:HB	1:C:110:ASP:HB3	1.99	0.43
1:G:75:TYR:CZ	2:G:201:BWU:N3	2.86	0.43
1:B:3:ILE:CD1	1:B:155:LEU:HB3	2.47	0.43
1:G:47:ASN:ND2	1:G:106:GLY:HA2	2.34	0.43
1:F:160:ASP:HB3	1:F:163:LEU:HD12	2.01	0.43
1:H:85:LEU:HD23	1:H:133:VAL:HG12	2.00	0.43
1:F:85:LEU:HD23	1:F:133:VAL:HG12	2.00	0.43
1:A:5:ILE:O	1:A:178:ALA:HA	2.18	0.43
1:E:130:ASP:O	1:E:135:LYS:HE3	2.17	0.43
1:D:11:TYR:OH	1:H:130:ASP:OD2	2.30	0.43
1:E:74:HIS:O	1:E:85:LEU:HA	2.19	0.42
1:H:5:ILE:O	1:H:178:ALA:HA	2.20	0.42
1:B:8:SER:HA	1:B:181:THR:O	2.19	0.42
1:E:139:GLY:CA	1:E:144:ASP:O	2.67	0.42
1:F:121:LEU:HD11	1:F:151:GLN:HG2	2.01	0.42
1:D:168:THR:HG23	1:D:171:GLN:H	1.83	0.42
1:A:84:THR:HG21	1:A:128:GLY:HA2	2.01	0.42
1:D:32:HIS:CD2	2:D:201:BWU:N4	2.87	0.42
1:D:38:VAL:HG21	1:D:42:ASN:HD21	1.84	0.42
1:D:130:ASP:OD2	1:F:59:LYS:NZ	2.52	0.42
1:H:47:ASN:HA	1:H:48:PRO:HA	1.90	0.42
1:C:79:SER:O	1:C:82:SER:HA	2.20	0.41
1:F:39:ASP:OD1	1:F:40:GLY:N	2.50	0.41
1:H:75:TYR:CE1	2:H:201:BWU:C40	3.03	0.41
1:F:95:GLY:O	1:F:96:ASP:OD1	2.38	0.41
1:F:160:ASP:OD1	1:F:160:ASP:C	2.64	0.41
1:G:181:THR:HG23	1:G:182:PRO:HD2	2.02	0.41
1:D:11:TYR:HA	1:D:14:TRP:CD1	2.56	0.41
1:A:12:SER:HB3	1:A:170:ASP:OD2	2.21	0.41
1:E:43:THR:HG23	1:E:58:LEU:CD1	2.51	0.41
1:A:77:LEU:O	1:A:82:SER:HA	2.21	0.41
1:B:47:ASN:HB3	1:B:55:GLN:HB2	2.03	0.41
1:B:58:LEU:HB3	1:B:67:PHE:CE2	2.55	0.41
1:E:165:ILE:HD12	1:E:165:ILE:HA	1.88	0.41
1:D:91:SER:HA	1:D:116:PHE:O	2.21	0.40
1:G:54:SER:HA	1:G:71:GLY:O	2.21	0.40
1:F:12:SER:HB2	1:F:170:ASP:OD1	2.21	0.40
1:H:84:THR:HG21	1:H:128:GLY:HA2	2.03	0.40
1:A:173:ALA:HA	1:A:178:ALA:O	2.20	0.40
1:D:5:ILE:O	1:D:178:ALA:HA	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:17:ALA:HB2	1:F:165:ILE:HG23	2.03	0.40
1:H:20:LEU:HD23	1:H:20:LEU:HA	1.88	0.40
2:H:201:BWU:C10	2:H:201:BWU:C21	2.97	0.40
1:A:17:ALA:HB2	1:A:165:ILE:HG23	2.03	0.40
1:A:91:SER:HA	1:A:116:PHE:O	2.21	0.40
1:C:89:LEU:HD21	1:C:92:ILE:HD11	2.03	0.40
1:F:172:LEU:HD13	1:F:172:LEU:HA	1.89	0.40
1:H:58:LEU:HD22	1:H:141:MET:HG2	2.04	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	179/184 (97%)	172 (96%)	7 (4%)	0	100	100
1	B	179/184 (97%)	171 (96%)	8 (4%)	0	100	100
1	C	173/184 (94%)	168 (97%)	5 (3%)	0	100	100
1	D	179/184 (97%)	173 (97%)	6 (3%)	0	100	100
1	E	179/184 (97%)	170 (95%)	9 (5%)	0	100	100
1	F	179/184 (97%)	177 (99%)	2 (1%)	0	100	100
1	G	179/184 (97%)	172 (96%)	7 (4%)	0	100	100
1	H	179/184 (97%)	176 (98%)	3 (2%)	0	100	100
All	All	1426/1472 (97%)	1379 (97%)	47 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	138/139 (99%)	137 (99%)	1 (1%)	76	89
1	B	138/139 (99%)	132 (96%)	6 (4%)	26	51
1	C	135/139 (97%)	129 (96%)	6 (4%)	25	50
1	D	138/139 (99%)	131 (95%)	7 (5%)	21	43
1	E	138/139 (99%)	131 (95%)	7 (5%)	21	43
1	F	138/139 (99%)	132 (96%)	6 (4%)	26	51
1	G	138/139 (99%)	131 (95%)	7 (5%)	21	43
1	H	138/139 (99%)	131 (95%)	7 (5%)	21	43
All	All	1101/1112 (99%)	1054 (96%)	47 (4%)	26	51

All (47) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	78	PHE
1	B	3	ILE
1	B	9	THR
1	B	64	ASP
1	B	78	PHE
1	B	115	SER
1	B	119	LEU
1	C	8	SER
1	C	36	GLN
1	C	37	VAL
1	C	43	THR
1	C	85	LEU
1	C	181	THR
1	D	38	VAL
1	D	41	SER
1	D	43	THR
1	D	73	LEU
1	D	103	SER
1	D	110	ASP

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Mol	Chain	Res	Type
1	D	172	LEU
1	E	31	ASN
1	E	37	VAL
1	E	42	ASN
1	E	80	ASN
1	E	111	SER
1	E	112	GLN
1	E	119	LEU
1	F	24	SER
1	F	52	ASP
1	F	64	ASP
1	F	155	LEU
1	F	172	LEU
1	F	177	VAL
1	G	6	SER
1	G	31	ASN
1	G	38	VAL
1	G	96	ASP
1	G	119	LEU
1	G	130	ASP
1	G	181	THR
1	H	10	THR
1	H	103	SER
1	H	104	SER
1	H	142	SER
1	H	155	LEU
1	H	163	LEU
1	H	172	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	112	GLN
1	A	151	GLN
1	B	151	GLN
1	C	179	HIS
1	D	36	GLN
1	D	42	ASN
1	E	42	ASN
1	F	80	ASN
1	F	112	GLN
1	F	127	GLN

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Mol	Chain	Res	Type
1	F	179	HIS
1	G	74	HIS
1	H	47	ASN
1	H	127	GLN
1	H	166	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

19 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	$\# Z > 2$
2	BWU	F	201	1	60,60,60	1.78	16 (26%)	72,94,94	1.63	14 (19%)
4	PO4	G	202	-	4,4,4	0.72	0	6,6,6	0.46	0
3	GOL	G	204	-	5,5,5	0.14	0	5,5,5	0.44	0
3	GOL	B	202	-	5,5,5	0.11	0	5,5,5	0.27	0
3	GOL	D	202	-	5,5,5	0.10	0	5,5,5	0.28	0
3	GOL	G	203	-	5,5,5	0.13	0	5,5,5	0.35	0
3	GOL	H	202	-	5,5,5	0.10	0	5,5,5	0.26	0
4	PO4	E	202	-	4,4,4	0.76	0	6,6,6	0.49	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	BWU	C	201	1	60,60,60	1.75	13 (21%)	72,94,94	1.61	13 (18%)
3	GOL	A	202	-	5,5,5	0.11	0	5,5,5	0.28	0
2	BWU	E	201	1	60,60,60	1.79	15 (25%)	72,94,94	1.63	14 (19%)
2	BWU	D	201	1	60,60,60	1.78	14 (23%)	72,94,94	1.62	13 (18%)
2	BWU	B	201	1	60,60,60	1.78	14 (23%)	72,94,94	1.69	13 (18%)
2	BWU	A	201	1	60,60,60	1.76	16 (26%)	72,94,94	1.62	14 (19%)
3	GOL	A	203	-	5,5,5	0.10	0	5,5,5	0.24	0
2	BWU	G	201	1	60,60,60	1.81	18 (30%)	72,94,94	1.64	12 (16%)
2	BWU	H	201	1	60,60,60	1.75	17 (28%)	72,94,94	1.63	14 (19%)
3	GOL	E	203	-	5,5,5	0.12	0	5,5,5	0.30	0
3	GOL	B	203	-	5,5,5	0.13	0	5,5,5	0.31	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
2	BWU	C	201	1	1/1/5/5	0/16/72/72	0/4/12/12
3	GOL	G	204	-	-	0/4/4/4	-
3	GOL	B	202	-	-	0/4/4/4	-
3	GOL	D	202	-	-	0/4/4/4	-
3	GOL	G	203	-	-	0/4/4/4	-
3	GOL	H	202	-	-	0/4/4/4	-
3	GOL	A	202	-	-	1/4/4/4	-
2	BWU	E	201	1	1/1/5/5	0/16/72/72	0/4/12/12
2	BWU	D	201	1	1/1/5/5	0/16/72/72	0/4/12/12
2	BWU	B	201	1	1/1/5/5	0/16/72/72	0/4/12/12
2	BWU	A	201	1	1/1/5/5	0/16/72/72	0/4/12/12
2	BWU	F	201	1	1/1/5/5	0/16/72/72	0/4/12/12
3	GOL	A	203	-	-	1/4/4/4	-
2	BWU	G	201	1	1/1/5/5	0/16/72/72	0/4/12/12
2	BWU	H	201	1	1/1/5/5	0/16/72/72	0/4/12/12
3	GOL	E	203	-	-	0/4/4/4	-
3	GOL	B	203	-	-	1/4/4/4	-

All (123) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	201	BWU	C25-C33	6.10	1.51	1.40
2	D	201	BWU	C25-C33	5.69	1.50	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	201	BWU	C25-C33	5.52	1.50	1.40
2	G	201	BWU	C25-C33	5.51	1.50	1.40
2	B	201	BWU	C25-C33	5.47	1.49	1.40
2	F	201	BWU	C25-C33	5.41	1.49	1.40
2	A	201	BWU	C25-C33	5.27	1.49	1.40
2	H	201	BWU	C25-C33	5.17	1.49	1.40
2	D	201	BWU	C26-C39	5.11	1.50	1.40
2	E	201	BWU	C26-C39	4.90	1.49	1.40
2	C	201	BWU	C26-C39	4.76	1.49	1.40
2	G	201	BWU	C26-C39	4.70	1.49	1.40
2	F	201	BWU	C26-C39	4.50	1.49	1.40
2	A	201	BWU	C26-C39	4.38	1.48	1.40
2	H	201	BWU	C26-C39	4.32	1.48	1.40
2	B	201	BWU	C26-C39	4.03	1.48	1.40
2	B	201	BWU	C28-C43	3.61	1.50	1.42
2	A	201	BWU	C28-C43	3.56	1.50	1.42
2	D	201	BWU	C28-C43	3.47	1.49	1.42
2	B	201	BWU	CO-N4	-3.42	1.73	1.96
2	F	201	BWU	C28-C43	3.40	1.49	1.42
2	F	201	BWU	C25-C26	-3.37	1.33	1.47
2	D	201	BWU	C37-C40	-3.33	1.37	1.44
2	H	201	BWU	C25-C26	-3.32	1.33	1.47
2	B	201	BWU	C25-C26	-3.30	1.34	1.47
2	C	201	BWU	C28-C43	3.29	1.49	1.42
2	G	201	BWU	C42-N4	-3.24	1.33	1.39
2	A	201	BWU	C25-C26	-3.23	1.34	1.47
2	C	201	BWU	C25-C26	-3.22	1.34	1.47
2	G	201	BWU	C41-C42	-3.20	1.37	1.44
2	A	201	BWU	C41-C42	-3.20	1.37	1.44
2	A	201	BWU	C37-C40	-3.17	1.38	1.44
2	G	201	BWU	C30-C36	-3.17	1.37	1.47
2	G	201	BWU	C35-C36	-3.17	1.37	1.44
2	H	201	BWU	C28-C43	3.15	1.49	1.42
2	F	201	BWU	C42-N4	-3.14	1.34	1.39
2	H	201	BWU	C42-N4	-3.14	1.34	1.39
2	B	201	BWU	C41-C42	-3.14	1.37	1.44
2	E	201	BWU	C42-N4	-3.11	1.34	1.39
2	A	201	BWU	C42-N4	-3.11	1.34	1.39
2	H	201	BWU	C37-C40	-3.09	1.38	1.44
2	D	201	BWU	C25-C26	-3.08	1.34	1.47
2	C	201	BWU	C37-C40	-3.06	1.38	1.44
2	G	201	BWU	C25-C26	-3.05	1.34	1.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	201	BWU	C41-C42	-3.05	1.38	1.44
2	D	201	BWU	C42-N4	-3.05	1.34	1.39
2	B	201	BWU	C37-C40	-3.05	1.38	1.44
2	F	201	BWU	C35-C36	-3.04	1.38	1.44
2	G	201	BWU	C37-C40	-3.04	1.38	1.44
2	B	201	BWU	C42-N4	-3.01	1.34	1.39
2	F	201	BWU	C41-C42	-3.01	1.38	1.44
2	E	201	BWU	C41-C42	-3.00	1.38	1.44
2	H	201	BWU	C41-C42	-2.99	1.38	1.44
2	G	201	BWU	C28-C43	2.99	1.48	1.42
2	C	201	BWU	C42-N4	-2.98	1.34	1.39
2	F	201	BWU	C37-C40	-2.97	1.38	1.44
2	E	201	BWU	C25-C26	-2.97	1.35	1.47
2	H	201	BWU	C35-C36	-2.93	1.38	1.44
2	B	201	BWU	C35-C36	-2.93	1.38	1.44
2	E	201	BWU	C30-C36	-2.91	1.38	1.47
2	E	201	BWU	C37-C40	-2.90	1.38	1.44
2	C	201	BWU	C35-C36	-2.89	1.38	1.44
2	D	201	BWU	C41-C42	-2.87	1.38	1.44
2	E	201	BWU	C35-C36	-2.85	1.38	1.44
2	F	201	BWU	C27-C29	2.83	1.49	1.41
2	C	201	BWU	CO-N4	-2.79	1.77	1.96
2	D	201	BWU	C35-C36	-2.75	1.38	1.44
2	B	201	BWU	C27-C29	2.73	1.49	1.41
2	F	201	BWU	C30-C36	-2.72	1.39	1.47
2	E	201	BWU	C28-C43	2.70	1.48	1.42
2	A	201	BWU	C35-C36	-2.70	1.38	1.44
2	H	201	BWU	C30-C36	-2.69	1.39	1.47
2	E	201	BWU	CO-N3	-2.69	1.70	1.91
2	H	201	BWU	C27-C29	2.69	1.49	1.41
2	A	201	BWU	C27-C29	2.68	1.49	1.41
2	A	201	BWU	C30-C36	-2.67	1.39	1.47
2	B	201	BWU	C30-C36	-2.66	1.39	1.47
2	E	201	BWU	C27-C29	2.61	1.49	1.41
2	F	201	BWU	C33-N2	-2.60	1.35	1.39
2	D	201	BWU	C30-C36	-2.55	1.39	1.47
2	C	201	BWU	C30-C36	-2.53	1.39	1.47
2	G	201	BWU	C27-C29	2.51	1.48	1.41
2	G	201	BWU	C36-N2	-2.48	1.33	1.38
2	H	201	BWU	C33-N2	-2.47	1.35	1.39
2	B	201	BWU	C33-N2	-2.43	1.35	1.39
2	G	201	BWU	CO-N4	-2.42	1.79	1.96

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	201	BWU	C27-C29	2.39	1.48	1.41
2	G	201	BWU	C30-N1	-2.39	1.34	1.38
2	B	201	BWU	C36-N2	-2.39	1.33	1.38
2	B	201	BWU	C38-C39	-2.38	1.39	1.44
2	D	201	BWU	C27-C29	2.38	1.48	1.41
2	A	201	BWU	C38-C39	-2.36	1.39	1.44
2	E	201	BWU	C36-N2	-2.34	1.33	1.38
2	D	201	BWU	C36-N2	-2.34	1.33	1.38
2	C	201	BWU	C36-N2	-2.31	1.33	1.38
2	F	201	BWU	CO-N4	-2.31	1.80	1.96
2	D	201	BWU	CO-N2	-2.31	1.80	1.96
2	A	201	BWU	CO-N3	-2.31	1.73	1.91
2	H	201	BWU	C36-N2	-2.31	1.33	1.38
2	G	201	BWU	C44-C43	-2.28	1.39	1.44
2	F	201	BWU	C36-N2	-2.27	1.34	1.38
2	A	201	BWU	C33-N2	-2.27	1.35	1.39
2	G	201	BWU	C33-N2	-2.24	1.35	1.39
2	C	201	BWU	C33-N2	-2.24	1.35	1.39
2	D	201	BWU	C33-N2	-2.22	1.35	1.39
2	E	201	BWU	C44-C43	-2.17	1.39	1.44
2	F	201	BWU	C30-N1	-2.17	1.35	1.38
2	H	201	BWU	C31-C30	-2.16	1.37	1.42
2	H	201	BWU	C44-C43	-2.15	1.39	1.44
2	H	201	BWU	CO-N4	-2.14	1.81	1.96
2	D	201	BWU	CO-N3	-2.09	1.75	1.91
2	A	201	BWU	C44-C43	-2.09	1.40	1.44
2	A	201	BWU	C40-N3	-2.08	1.35	1.38
2	A	201	BWU	C36-N2	-2.08	1.34	1.38
2	E	201	BWU	C30-N1	-2.07	1.35	1.38
2	H	201	BWU	C38-C39	-2.07	1.40	1.44
2	F	201	BWU	C31-C30	-2.07	1.37	1.42
2	E	201	BWU	C33-N2	-2.05	1.35	1.39
2	G	201	BWU	C31-C30	-2.04	1.38	1.42
2	G	201	BWU	C40-N3	-2.04	1.35	1.38
2	F	201	BWU	C44-C43	-2.04	1.40	1.44
2	G	201	BWU	C38-C39	-2.03	1.40	1.44
2	H	201	BWU	C34-C33	-2.01	1.40	1.44

All (107) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	201	BWU	C35-C36-C30	-6.77	123.75	139.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	201	BWU	C35-C36-C30	-6.74	123.84	139.41
2	G	201	BWU	C35-C36-C30	-6.73	123.85	139.41
2	F	201	BWU	C35-C36-C30	-6.71	123.90	139.41
2	H	201	BWU	C35-C36-C30	-6.63	124.09	139.41
2	A	201	BWU	C35-C36-C30	-6.62	124.11	139.41
2	E	201	BWU	C35-C36-C30	-6.56	124.24	139.41
2	D	201	BWU	C35-C36-C30	-6.53	124.31	139.41
2	B	201	BWU	C31-C30-C36	-6.29	123.30	139.72
2	H	201	BWU	C31-C30-C36	-6.22	123.48	139.72
2	A	201	BWU	C31-C30-C36	-6.19	123.55	139.72
2	F	201	BWU	C31-C30-C36	-6.17	123.61	139.72
2	G	201	BWU	C31-C30-C36	-6.14	123.69	139.72
2	E	201	BWU	C31-C30-C36	-6.08	123.84	139.72
2	D	201	BWU	C31-C30-C36	-6.05	123.92	139.72
2	C	201	BWU	C31-C30-C36	-5.97	124.13	139.72
2	E	201	BWU	C41-C42-C40	-3.20	123.64	137.19
2	C	201	BWU	C41-C42-C40	-3.18	123.73	137.19
2	D	201	BWU	C41-C42-C40	-3.16	123.81	137.19
2	F	201	BWU	C26-C25-C33	-3.13	124.48	129.47
2	B	201	BWU	C41-C42-C40	-3.13	123.96	137.19
2	F	201	BWU	C25-C26-C39	-3.11	124.64	129.24
2	H	201	BWU	C41-C42-C40	-3.11	124.05	137.19
2	B	201	BWU	C26-C25-C33	-3.09	124.55	129.47
2	H	201	BWU	C25-C26-C39	-3.08	124.69	129.24
2	G	201	BWU	C41-C42-C40	-3.08	124.17	137.19
2	F	201	BWU	C41-C42-C40	-3.07	124.21	137.19
2	C	201	BWU	C25-C26-C39	-3.06	124.72	129.24
2	D	201	BWU	C26-C25-C33	-3.05	124.61	129.47
2	A	201	BWU	C41-C42-C40	-3.05	124.30	137.19
2	H	201	BWU	C26-C25-C33	-3.04	124.63	129.47
2	D	201	BWU	C41-C42-N4	3.01	113.21	109.65
2	B	201	BWU	C25-C26-C39	-2.98	124.83	129.24
2	D	201	BWU	C44-C41-C42	-2.95	103.99	107.42
2	A	201	BWU	C26-C25-C33	-2.94	124.80	129.47
2	B	201	BWU	C32-C31-C30	-2.91	103.83	107.47
2	E	201	BWU	C34-C35-C36	-2.89	104.06	107.42
2	C	201	BWU	C26-C25-C33	-2.86	124.92	129.47
2	B	201	BWU	C34-C35-C36	-2.80	104.16	107.42
2	G	201	BWU	C34-C35-C36	-2.79	104.17	107.42
2	F	201	BWU	C37-C40-C42	-2.77	123.36	136.50
2	B	201	BWU	C44-C41-C42	-2.75	104.22	107.42
2	A	201	BWU	C25-C26-C39	-2.75	125.17	129.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	201	BWU	C25-C26-C39	-2.74	125.19	129.24
2	H	201	BWU	C37-C40-C42	-2.72	123.59	136.50
2	C	201	BWU	C38-C37-C40	-2.72	104.31	107.55
2	G	201	BWU	C25-C26-C39	-2.71	125.24	129.24
2	D	201	BWU	C37-C40-C42	-2.70	123.68	136.50
2	C	201	BWU	C37-C40-C42	-2.68	123.80	136.50
2	F	201	BWU	C34-C35-C36	-2.67	104.31	107.42
2	B	201	BWU	C31-C30-N1	2.67	113.17	109.89
2	B	201	BWU	C37-C40-C42	-2.66	123.87	136.50
2	C	201	BWU	C41-C42-N4	2.66	112.79	109.65
2	C	201	BWU	C44-C41-C42	-2.66	104.33	107.42
2	A	201	BWU	C37-C40-C42	-2.65	123.91	136.50
2	H	201	BWU	C34-C35-C36	-2.63	104.36	107.42
2	E	201	BWU	C37-C40-C42	-2.62	124.09	136.50
2	E	201	BWU	C38-C37-C40	-2.61	104.44	107.55
2	B	201	BWU	C41-C42-N4	2.61	112.74	109.65
2	G	201	BWU	C26-C25-C33	-2.60	125.33	129.47
2	F	201	BWU	C38-C37-C40	-2.58	104.47	107.55
2	D	201	BWU	C38-C37-C40	-2.58	104.47	107.55
2	E	201	BWU	C25-C26-C39	-2.58	125.43	129.24
2	G	201	BWU	C37-C40-C42	-2.55	124.42	136.50
2	C	201	BWU	C34-C35-C36	-2.53	104.47	107.42
2	A	201	BWU	C34-C35-C36	-2.53	104.47	107.42
2	E	201	BWU	C26-C25-C33	-2.51	125.47	129.47
2	A	201	BWU	C32-C31-C30	-2.50	104.33	107.47
2	G	201	BWU	C38-C37-C40	-2.48	104.59	107.55
2	D	201	BWU	C34-C35-C36	-2.48	104.53	107.42
2	E	201	BWU	C35-C36-N2	2.46	113.40	110.18
2	G	201	BWU	C32-C31-C30	-2.42	104.43	107.47
2	E	201	BWU	C41-C42-N4	2.39	112.47	109.65
2	G	201	BWU	C35-C36-N2	2.36	113.26	110.18
2	H	201	BWU	C41-C42-N4	2.34	112.42	109.65
2	H	201	BWU	C32-C31-C30	-2.34	104.54	107.47
2	A	201	BWU	C38-C37-C40	-2.34	104.77	107.55
2	H	201	BWU	C27-C28-C43	-2.33	124.20	130.25
2	G	201	BWU	C31-C30-N1	2.31	112.73	109.89
2	H	201	BWU	C38-C37-C40	-2.31	104.80	107.55
2	F	201	BWU	C44-C41-C42	-2.31	104.73	107.42
2	A	201	BWU	C31-C30-N1	2.30	112.71	109.89
2	A	201	BWU	C27-C28-C43	-2.30	124.30	130.25
2	F	201	BWU	C41-C42-N4	2.29	112.36	109.65
2	E	201	BWU	C44-C41-C42	-2.28	104.77	107.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	201	BWU	C32-C31-C30	-2.28	104.62	107.47
2	F	201	BWU	C27-C28-C43	-2.26	124.39	130.25
2	A	201	BWU	C44-C41-C42	-2.25	104.80	107.42
2	B	201	BWU	C27-C28-C43	-2.24	124.45	130.25
2	B	201	BWU	C35-C36-N2	2.20	113.05	110.18
2	E	201	BWU	C32-C31-C30	-2.20	104.71	107.47
2	H	201	BWU	C44-C41-C42	-2.19	104.87	107.42
2	E	201	BWU	C27-C28-C43	-2.17	124.63	130.25
2	D	201	BWU	C27-C28-C43	-2.16	124.65	130.25
2	F	201	BWU	C32-C31-C30	-2.15	104.78	107.47
2	G	201	BWU	C27-C28-C43	-2.13	124.73	130.25
2	A	201	BWU	C41-C42-N4	2.11	112.15	109.65
2	F	201	BWU	C35-C36-N2	2.08	112.90	110.18
2	C	201	BWU	C31-C30-N1	2.08	112.44	109.89
2	D	201	BWU	C35-C36-N2	2.06	112.87	110.18
2	A	201	BWU	C35-C36-N2	2.05	112.86	110.18
2	F	201	BWU	C37-C40-N3	2.04	112.64	109.95
2	H	201	BWU	C31-C30-N1	2.04	112.40	109.89
2	D	201	BWU	C37-C40-N3	2.03	112.62	109.95
2	E	201	BWU	C31-C30-N1	2.03	112.39	109.89
2	H	201	BWU	C35-C36-N2	2.01	112.81	110.18
2	C	201	BWU	C37-C40-N3	2.01	112.59	109.95

All (8) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	A	201	BWU	N3
2	B	201	BWU	N3
2	C	201	BWU	N3
2	D	201	BWU	N3
2	E	201	BWU	N3
2	F	201	BWU	N3
2	G	201	BWU	N3
2	H	201	BWU	N3

All (3) torsion outliers are listed below:

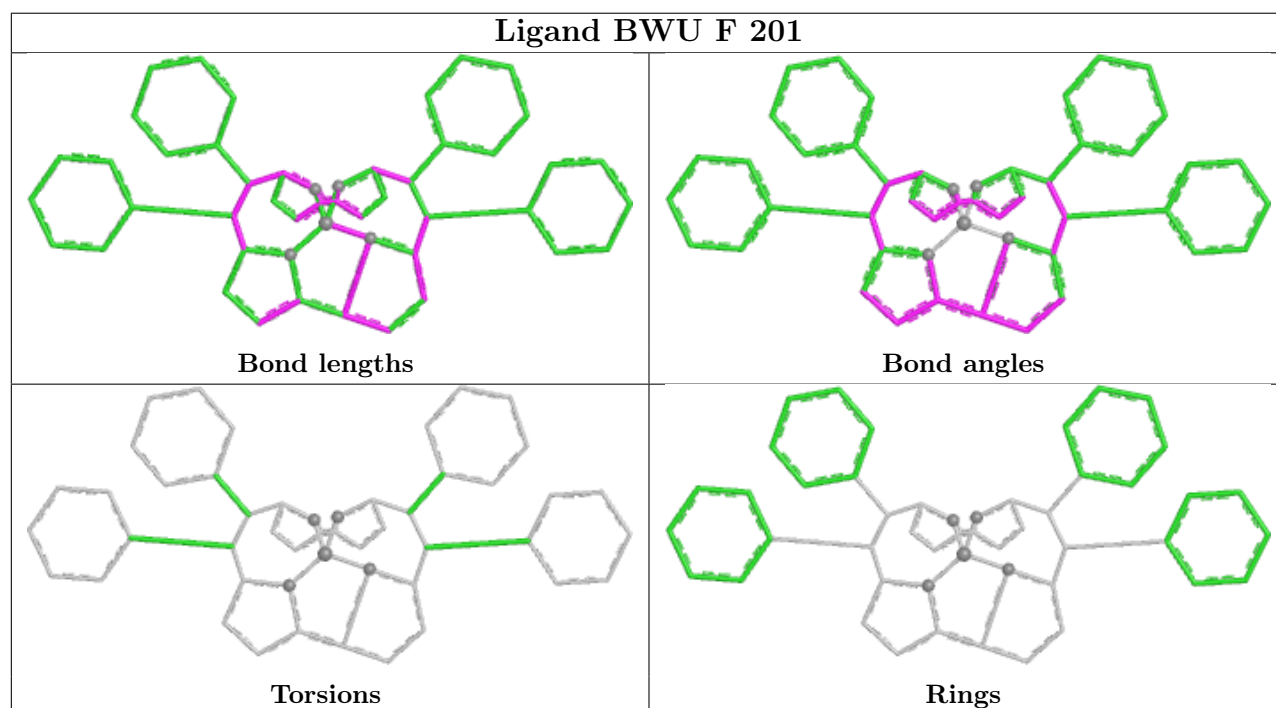
Mol	Chain	Res	Type	Atoms
3	A	203	GOL	O2-C2-C3-O3
3	B	203	GOL	O1-C1-C2-O2
3	A	202	GOL	O2-C2-C3-O3

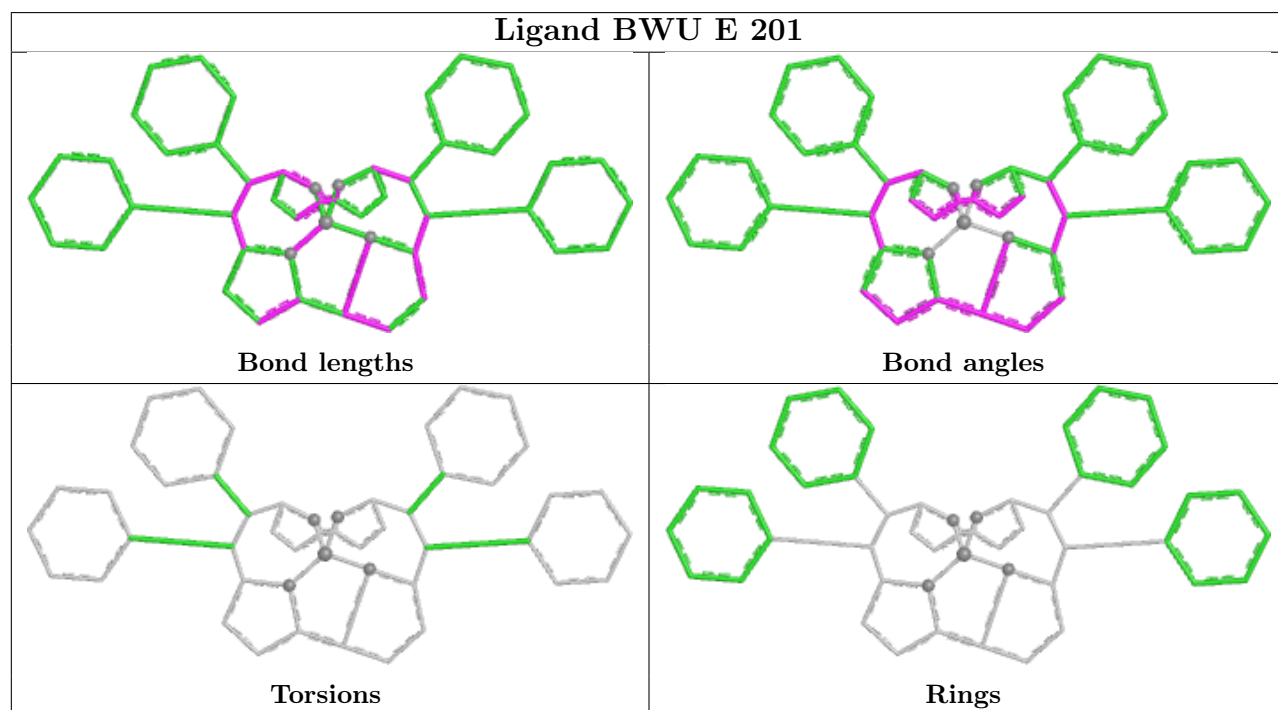
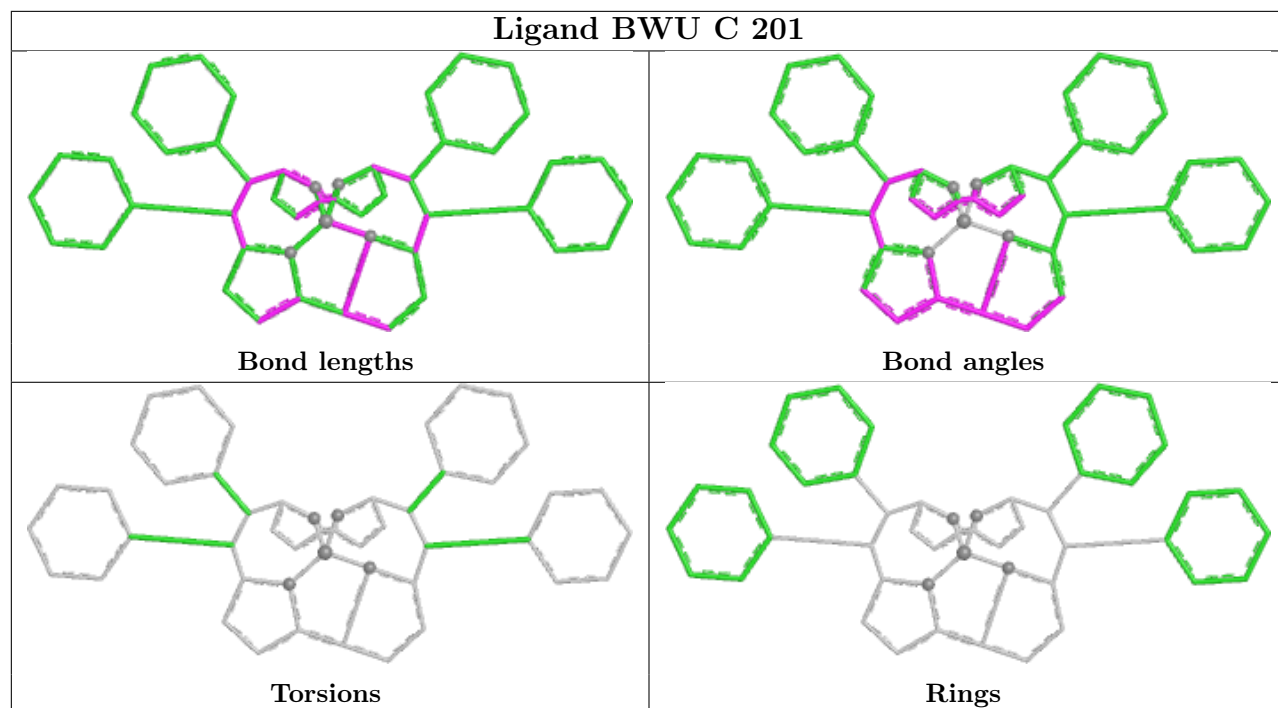
There are no ring outliers.

7 monomers are involved in 9 short contacts:

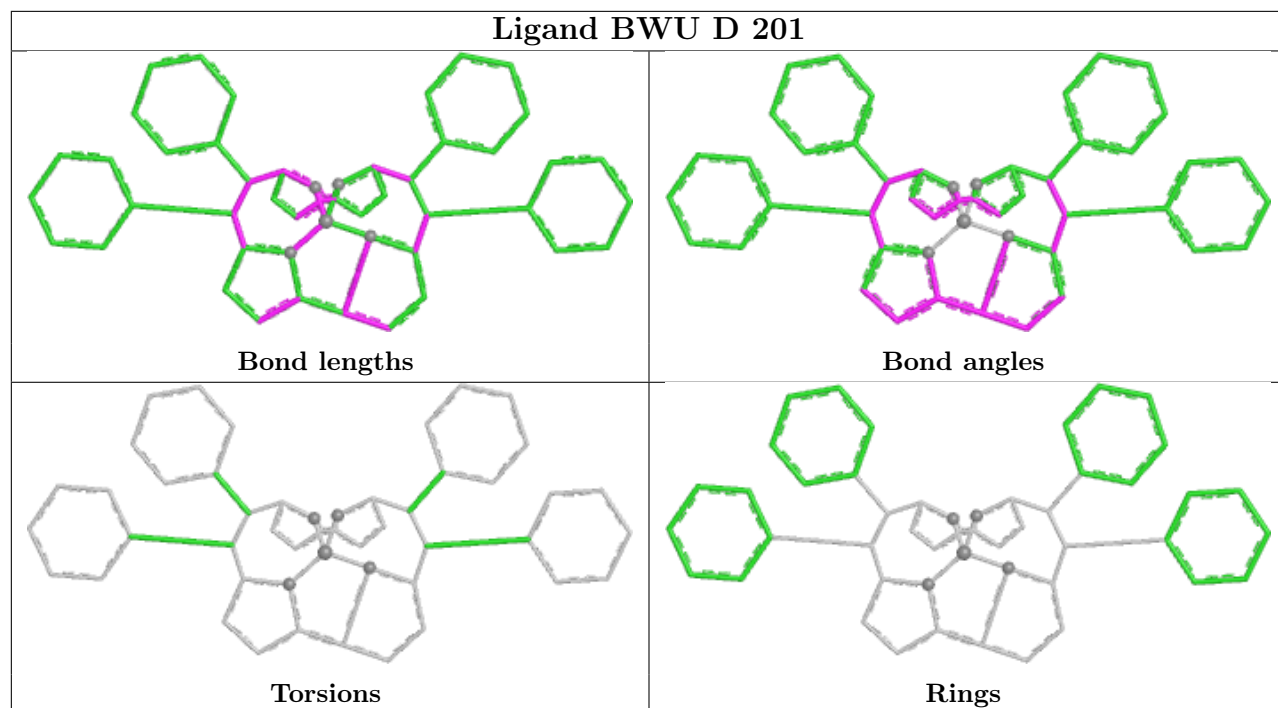
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	G	202	PO4	1	0
3	G	204	GOL	1	0
3	D	202	GOL	2	0
3	A	202	GOL	1	0
2	D	201	BWU	1	0
2	G	201	BWU	1	0
2	H	201	BWU	2	0

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.

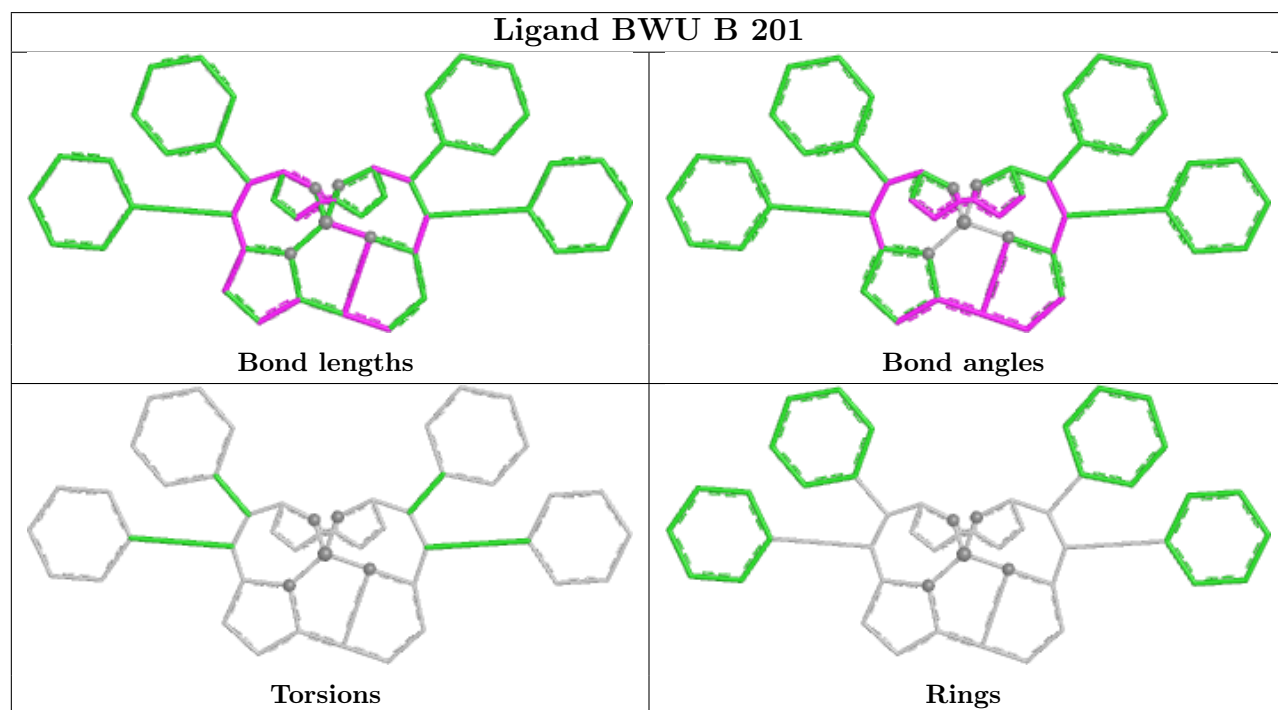


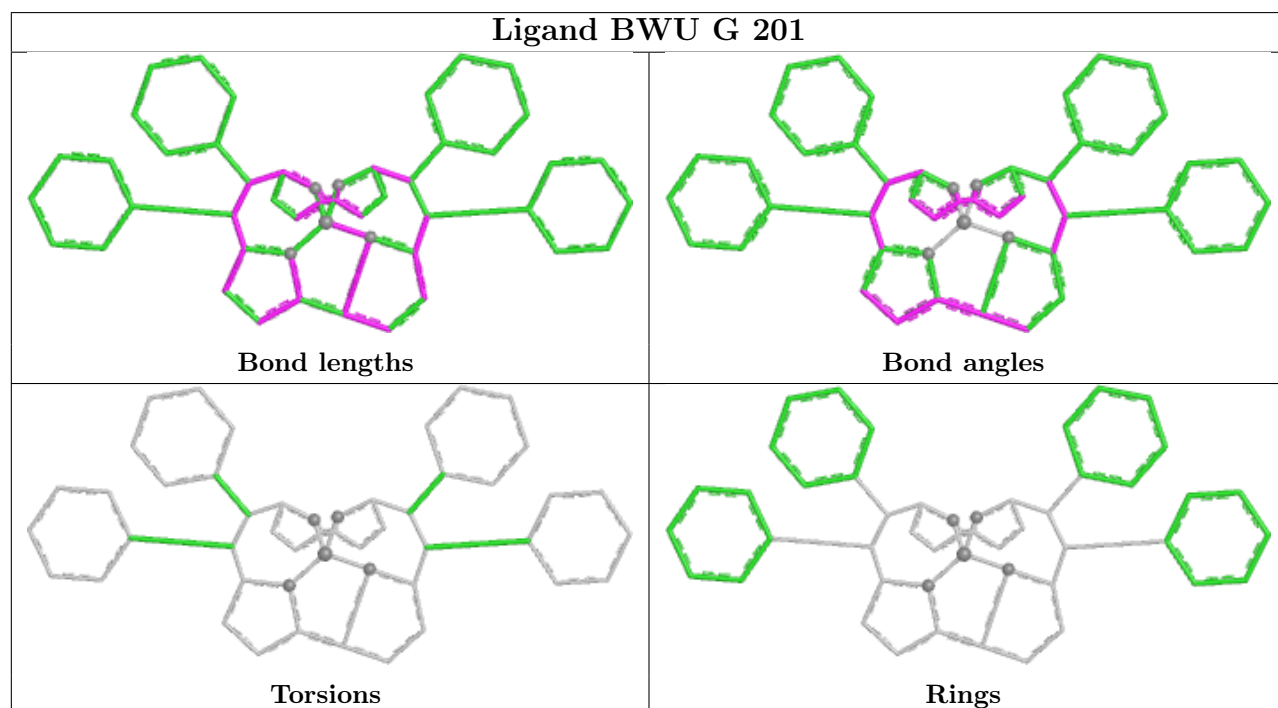
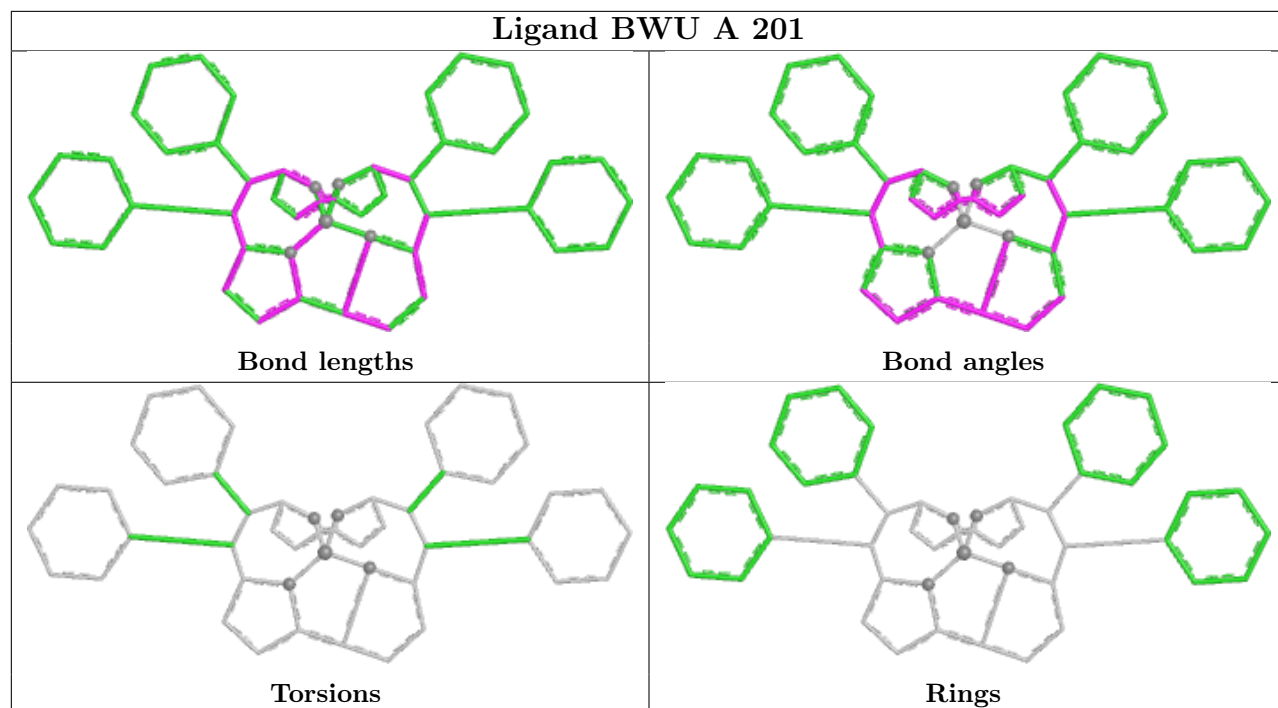


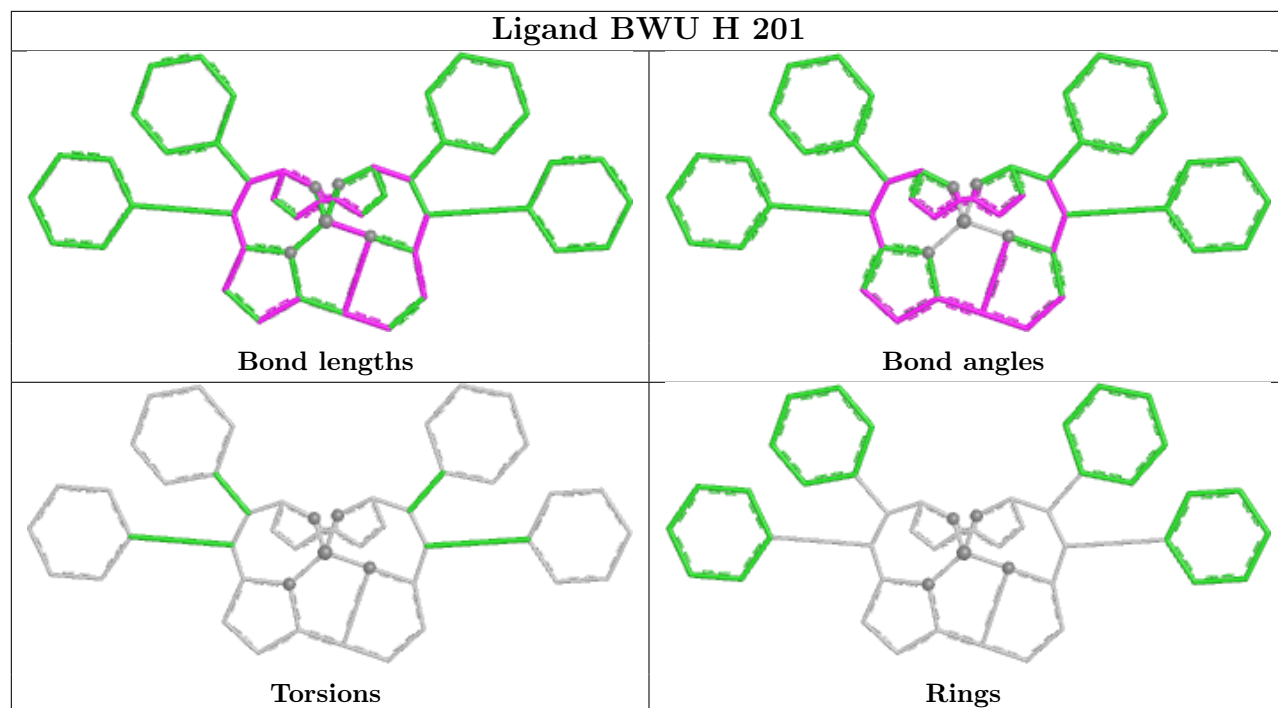
Ligand BWU D 201



Ligand BWU B 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 ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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	181/184 (98%)	0.63	19 (10%)	11 10	28, 45, 81, 117	0
1	B	181/184 (98%)	0.53	17 (9%)	14 12	28, 44, 84, 112	0
1	C	177/184 (96%)	0.55	13 (7%)	21 19	34, 55, 86, 95	0
1	D	181/184 (98%)	0.72	20 (11%)	10 8	31, 55, 92, 123	0
1	E	181/184 (98%)	0.30	9 (4%)	34 30	27, 40, 63, 132	0
1	F	181/184 (98%)	1.17	35 (19%)	3 2	39, 66, 110, 126	0
1	G	181/184 (98%)	0.41	11 (6%)	27 23	28, 41, 65, 136	0
1	H	181/184 (98%)	0.82	16 (8%)	15 14	37, 60, 93, 112	0
All	All	1444/1472 (98%)	0.64	140 (9%)	13 11	27, 50, 94, 136	0

All (140) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	38	VAL	6.8
1	D	38	VAL	6.1
1	E	37	VAL	5.7
1	C	38	VAL	5.5
1	D	39	ASP	5.4
1	E	38	VAL	5.2
1	A	110	ASP	5.1
1	G	40	GLY	5.0
1	D	41	SER	5.0
1	C	43	THR	4.9
1	D	40	GLY	4.8
1	B	160	ASP	4.7
1	D	37	VAL	4.7
1	C	37	VAL	4.6
1	G	37	VAL	4.6
1	A	95	GLY	4.6

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Mol	Chain	Res	Type	RSRZ
1	E	43	THR	4.4
1	H	163	LEU	4.4
1	E	40	GLY	4.4
1	B	182	PRO	4.3
1	A	182	PRO	4.3
1	A	77	LEU	4.3
1	G	43	THR	4.0
1	D	42	ASN	4.0
1	A	161	PRO	4.0
1	F	105	GLY	3.8
1	F	103	SER	3.8
1	H	182	PRO	3.8
1	H	165	ILE	3.8
1	F	177	VAL	3.7
1	D	44	GLY	3.7
1	C	102	ALA	3.7
1	B	78	PHE	3.7
1	B	110	ASP	3.7
1	B	161	PRO	3.6
1	F	182	PRO	3.6
1	H	104	SER	3.5
1	F	180	ALA	3.4
1	G	41	SER	3.4
1	B	77	LEU	3.4
1	F	102	ALA	3.4
1	G	39	ASP	3.4
1	E	39	ASP	3.3
1	F	13	GLY	3.3
1	F	77	LEU	3.2
1	A	97	THR	3.2
1	A	3	ILE	3.2
1	H	102	ALA	3.2
1	D	163	LEU	3.2
1	F	18	ASP	3.1
1	F	3	ILE	3.1
1	D	36	GLN	3.1
1	F	174	ALA	3.0
1	F	179	HIS	3.0
1	B	177	VAL	3.0
1	D	43	THR	3.0
1	E	179	HIS	3.0
1	F	6	SER	3.0

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Mol	Chain	Res	Type	RSRZ
1	A	122	ASP	2.9
1	F	111	SER	2.9
1	F	161	PRO	2.9
1	F	169	PHE	2.9
1	H	72	ASP	2.8
1	A	159	VAL	2.8
1	H	105	GLY	2.8
1	A	158	ALA	2.8
1	C	72	ASP	2.7
1	F	104	SER	2.7
1	G	126	ALA	2.7
1	A	179	HIS	2.7
1	E	182	PRO	2.7
1	D	104	SER	2.7
1	F	163	LEU	2.7
1	B	176	GLY	2.6
1	E	42	ASN	2.6
1	F	19	TYR	2.6
1	F	97	THR	2.6
1	D	182	PRO	2.6
1	H	96	ASP	2.6
1	C	163	LEU	2.6
1	B	175	ALA	2.6
1	C	158	ALA	2.6
1	B	170	ASP	2.5
1	H	101	GLY	2.5
1	B	179	HIS	2.5
1	A	76	THR	2.5
1	H	181	THR	2.5
1	D	110	ASP	2.5
1	G	42	ASN	2.5
1	C	2	SER	2.5
1	E	41	SER	2.5
1	F	165	ILE	2.4
1	F	79	SER	2.4
1	C	85	LEU	2.4
1	D	102	ALA	2.4
1	H	10	THR	2.3
1	A	162	SER	2.3
1	B	3	ILE	2.3
1	C	110	ASP	2.3
1	H	106	GLY	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	90	ASP	2.3
1	B	9	THR	2.3
1	A	2	SER	2.2
1	F	5	ILE	2.2
1	D	52	ASP	2.2
1	B	76	THR	2.2
1	B	181	THR	2.2
1	H	100	GLY	2.2
1	H	162	SER	2.2
1	A	177	VAL	2.2
1	G	88	LYS	2.2
1	F	181	THR	2.2
1	D	165	ILE	2.2
1	A	160	ASP	2.2
1	F	16	VAL	2.2
1	A	175	ALA	2.2
1	F	170	ASP	2.2
1	H	64	ASP	2.2
1	F	14	TRP	2.2
1	G	80	ASN	2.1
1	D	170	ASP	2.1
1	F	110	ASP	2.1
1	F	159	VAL	2.1
1	F	80	ASN	2.1
1	F	106	GLY	2.1
1	G	125	ILE	2.1
1	F	122	ASP	2.1
1	D	72	ASP	2.1
1	C	36	GLN	2.1
1	F	116	PHE	2.1
1	F	119	LEU	2.1
1	D	2	SER	2.1
1	A	78	PHE	2.0
1	B	158	ALA	2.0
1	F	98	LEU	2.0
1	H	14	TRP	2.0
1	C	182	PRO	2.0
1	D	103	SER	2.0
1	B	164	SER	2.0
1	C	103	SER	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

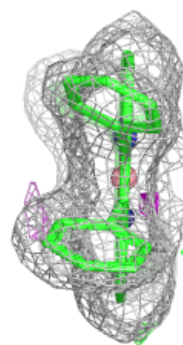
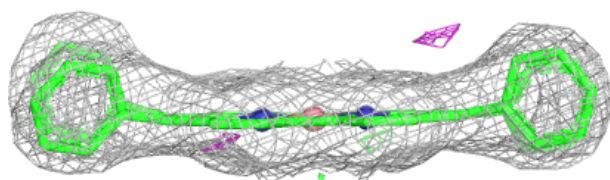
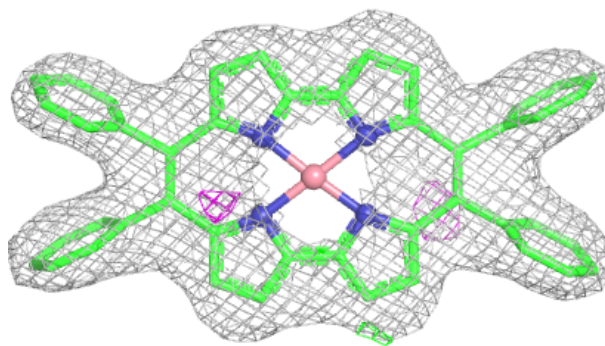
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	GOL	A	202	6/6	0.73	0.14	51,54,55,56	0
3	GOL	A	203	6/6	0.76	0.14	48,49,52,53	0
3	GOL	B	203	6/6	0.77	0.19	40,48,50,52	0
3	GOL	G	203	6/6	0.78	0.16	42,52,56,56	0
3	GOL	B	202	6/6	0.82	0.18	45,57,58,58	0
3	GOL	E	203	6/6	0.88	0.14	50,55,56,57	0
3	GOL	G	204	6/6	0.88	0.14	39,47,48,50	0
3	GOL	H	202	6/6	0.89	0.15	51,55,55,58	0
3	GOL	D	202	6/6	0.95	0.12	41,47,50,50	0
2	BWU	F	201	49/49	0.96	0.09	37,40,49,51	0
2	BWU	B	201	49/49	0.96	0.09	26,30,34,36	0
2	BWU	A	201	49/49	0.97	0.08	27,31,34,35	0
2	BWU	G	201	49/49	0.97	0.08	26,31,39,41	0
2	BWU	H	201	49/49	0.97	0.09	36,40,45,46	0
2	BWU	E	201	49/49	0.97	0.07	24,28,38,39	0
4	PO4	E	202	5/5	0.97	0.10	34,41,45,55	0
4	PO4	G	202	5/5	0.97	0.08	35,41,46,53	0
2	BWU	D	201	49/49	0.98	0.08	28,31,51,57	0
2	BWU	C	201	49/49	0.98	0.08	30,35,50,56	0

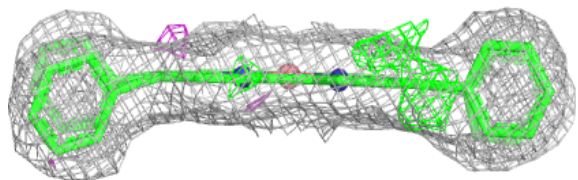
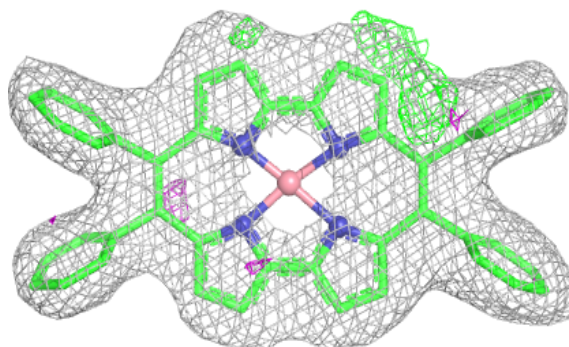
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around BWU F 201:

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

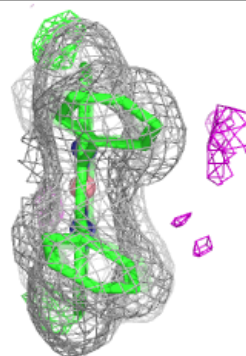
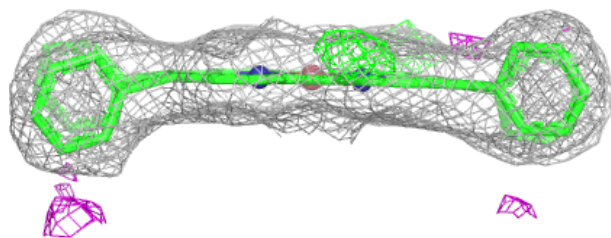
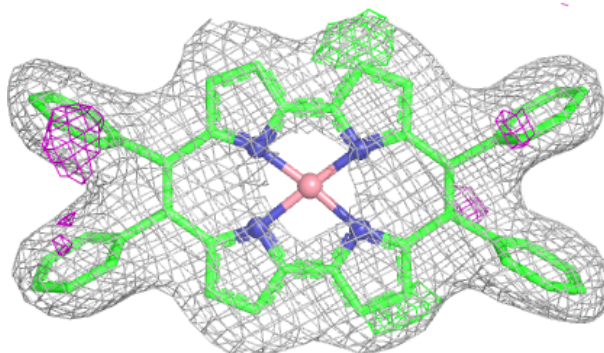
**Electron density around BWU B 201:**

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

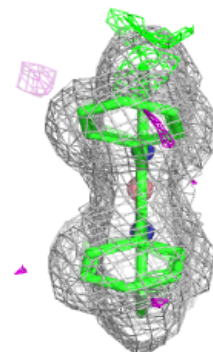
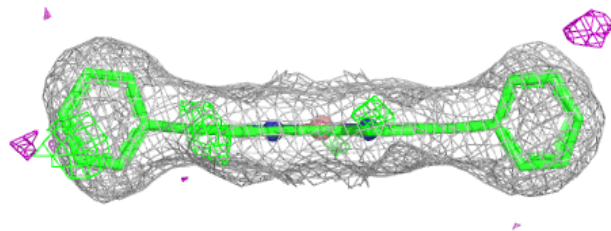
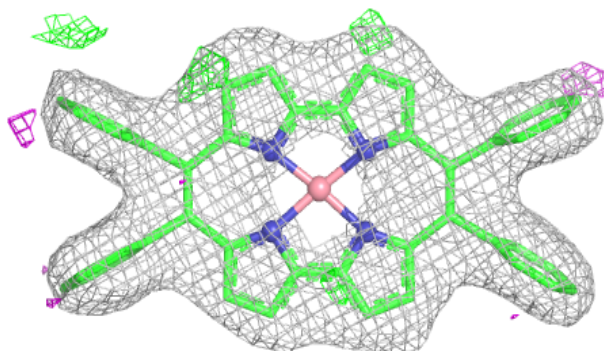


Electron density around BWU A 201:

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

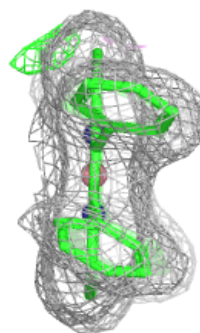
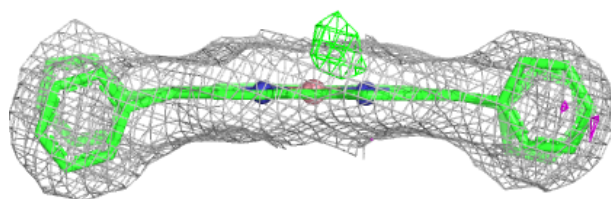
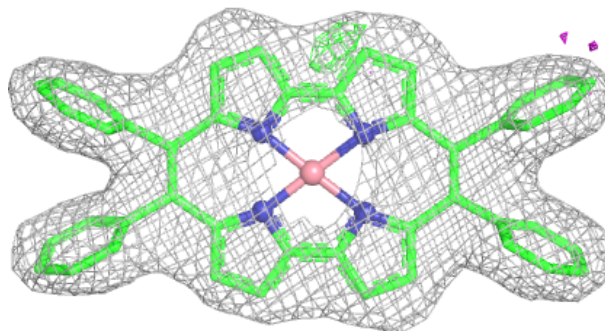
**Electron density around BWU G 201:**

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

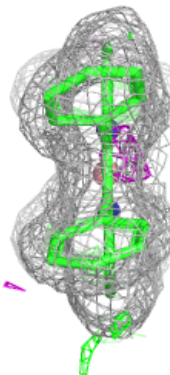
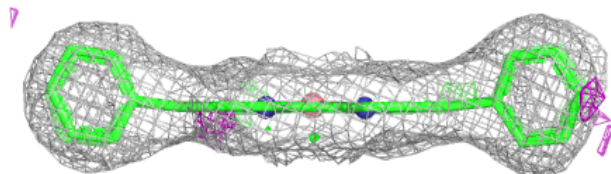
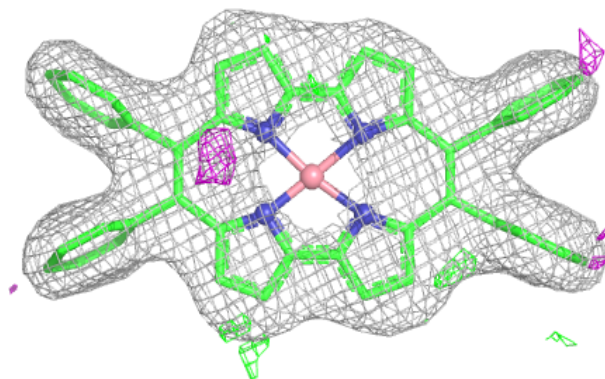


Electron density around BWU H 201:

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

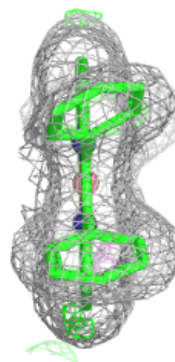
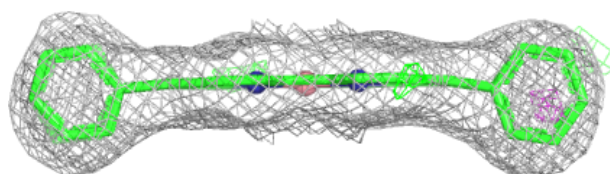
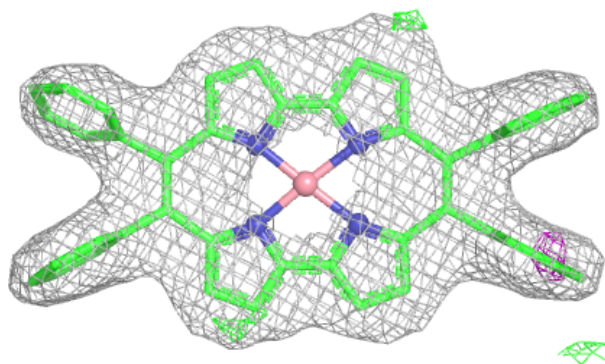
**Electron density around BWU E 201:**

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

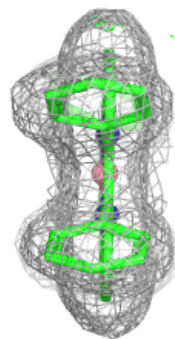
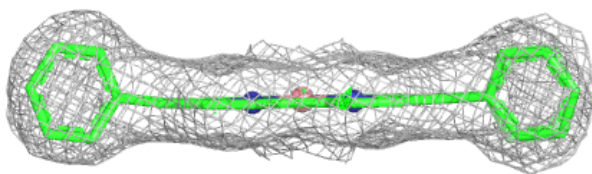
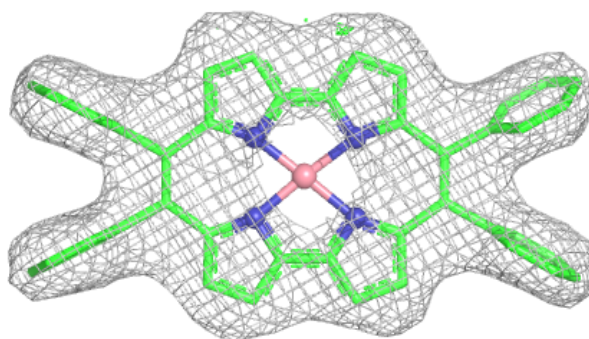


Electron density around BWU D 201:

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

**Electron density around BWU 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)



6.5 Other polymers [i](#)

There are no such residues in this entry.