



Full wwPDB EM Validation Report ⓘ

Aug 5, 2026 – 01:36 AM EDT

PDB ID : 8EUN / pdb_00008eun
EMDB ID : EMD-28616
Title : MicroED structure of an Aeropyrum pernix protoglobin metallo-carbene complex
Authors : Danelius, E.; Gonen, T.; Unge, J.T.
Deposited on : 2022-10-19
Resolution : 2.50 Å(reported)

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

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/EMValidationReportHelp>
with specific help available everywhere you see the ⓘ 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:

EMDB validation analysis	: NOT EXECUTED
Mogul	: 2022.3.0, CSD as543be (2022)
MolProbity	: 4-5-2 with Phenix2.0
Buster-report	: wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	: 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics	: NOT EXECUTED
MapQ	: NOT EXECUTED
CCP4	: 9.0.015 (Gargrove)
Density-Fitness	: 1.0.12
Ideal geometry (proteins)	: Engh & Huber (2001)
Ideal geometry (DNA, RNA)	: Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	: 2.50

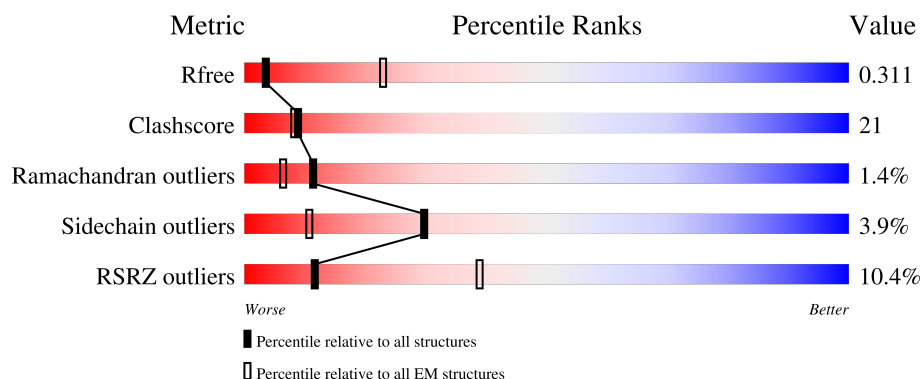
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON CRYSTALLOGRAPHY

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)	EM structures (#Entries)
R_{free}	180332	208
Clashscore	229148	23984
Ramachandran outliers	224038	23583
Sidechain outliers	223484	23102
RSRZ outliers	180361	209

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

Mol	Chain	Length	Quality of chain
1	A	195	11% 50% 44% • •
1	B	195	9% 62% 32% • •

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3268 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

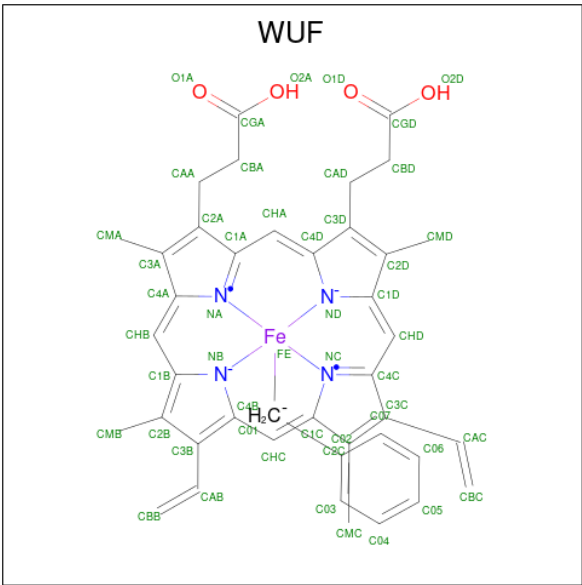
- Molecule 1 is a protein called Protogloblin ApPgb.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	187	Total	C	N	O	S	0	0
			1552	1008	260	281	3		
1	B	187	Total	C	N	O	S	0	0
			1552	1008	260	281	3		

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	45	GLY	CYS	engineered mutation	UNP Q9YFF4
A	59	LEU	TRP	engineered mutation	UNP Q9YFF4
A	60	VAL	TYR	engineered mutation	UNP Q9YFF4
A	63	ARG	VAL	engineered mutation	UNP Q9YFF4
A	102	SER	CYS	engineered mutation	UNP Q9YFF4
A	145	GLN	PHE	engineered mutation	UNP Q9YFF4
A	149	LEU	ILE	engineered mutation	UNP Q9YFF4
B	45	GLY	CYS	engineered mutation	UNP Q9YFF4
B	59	LEU	TRP	engineered mutation	UNP Q9YFF4
B	60	VAL	TYR	engineered mutation	UNP Q9YFF4
B	63	ARG	VAL	engineered mutation	UNP Q9YFF4
B	102	SER	CYS	engineered mutation	UNP Q9YFF4
B	145	GLN	PHE	engineered mutation	UNP Q9YFF4
B	149	LEU	ILE	engineered mutation	UNP Q9YFF4

- Molecule 2 is benzyl[3,3'-(7,12-diethenyl-3,8,13,17-tetramethylporphyrin-2,18-diyl-kappa 4 N 21 ,N 22 ,N 23 ,N 24)di(propanoato)(2-)]iron (CCD ID: WUF) (formula: C₄₁H₃₉FeN₄O₄) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					AltConf
2	A	1	Total	C	Fe	N	O	0
			50	41	1	4	4	
2	B	1	Total	C	Fe	N	O	0
			50	41	1	4	4	

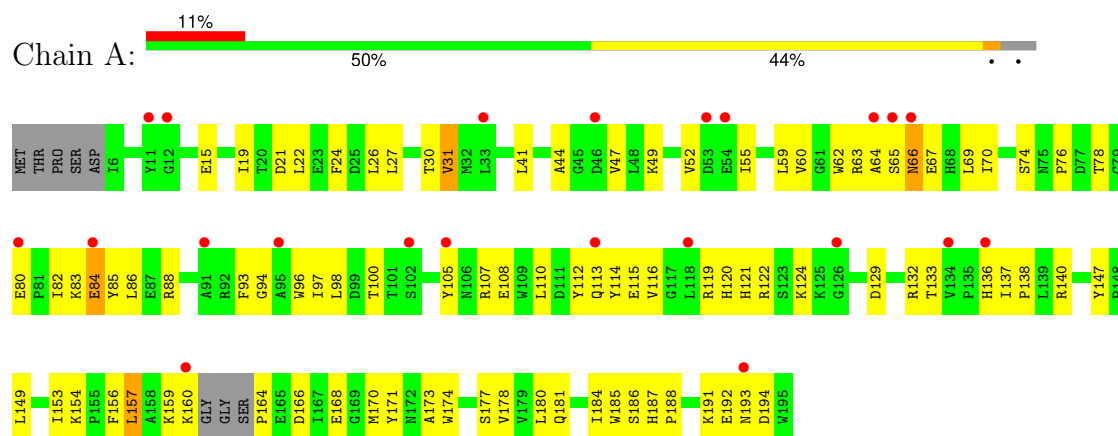
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		AltConf
3	A	33	Total	O	0
			33	33	
3	B	31	Total	O	0
			31	31	

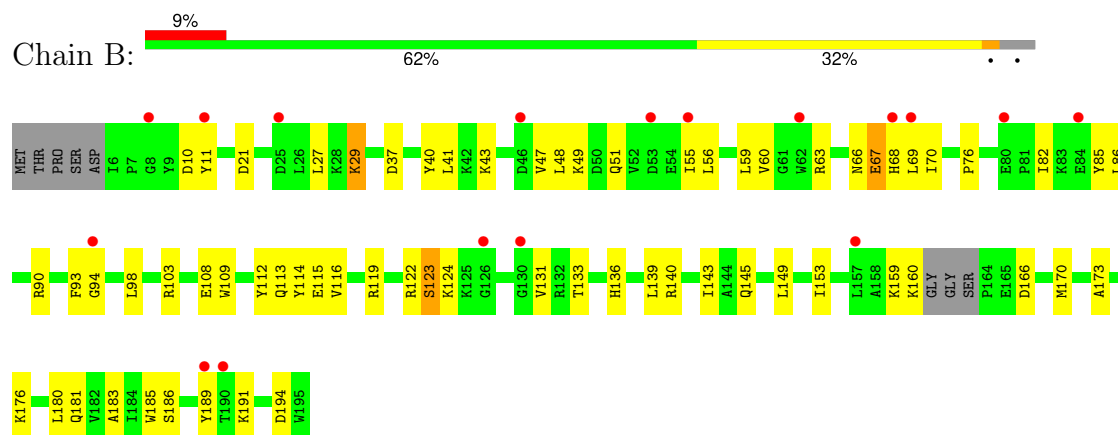
3 Residue-property plots [i](#)

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. 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. 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: Protogloblin ApPgb



• Molecule 1: Protogloblin ApPgb



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	58.12Å 45.76Å 72.63Å 90.00° 105.72° 90.00°	Depositor
Resolution (Å)	24.37 – 2.50 24.37 – 2.50	Depositor EDS
% Data completeness (in resolution range)	70.4 (24.37-2.50) 70.3 (24.37-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.95 (at 2.50Å)	Xtriage
Refinement program	phenix.refine 1.20.1_4487, PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.233 , 0.284 0.258 , 0.311	Depositor DCC
R_{free} test set	898 reflections (6.92%)	wwPDB-VP
Wilson B-factor (Å ²)	17.6	Xtriage
Anisotropy	1.449	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.41 , 47.7	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	0.84	EDS
Total number of atoms	3268	wwPDB-VP
Average B, all atoms (Å ²)	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 11.99% 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: WUF

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	0.18	0/1595	0.41	0/2163
1	B	0.20	0/1595	0.40	0/2163
All	All	0.19	0/3190	0.40	0/4326

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1552	0	1537	78	1
1	B	1552	0	1537	56	0
2	A	50	0	0	2	0
2	B	50	0	0	4	0
3	A	33	0	0	20	1
3	B	31	0	0	13	0
All	All	3268	0	3074	131	2

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 21.

All (131) 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:180:LEU:O	3:B:301:HOH:O	1.77	1.02
1:B:59:LEU:O	1:B:66:ASN:ND2	2.09	0.85
1:A:191:LYS:N	3:B:302:HOH:O	2.10	0.82
1:A:62:TRP:CD1	1:A:63:ARG:HH21	1.99	0.80
1:A:110:LEU:O	3:A:301:HOH:O	1.99	0.79
1:B:113:GLN:NE2	3:B:303:HOH:O	2.17	0.77
1:A:96:TRP:NE1	3:A:302:HOH:O	2.18	0.76
1:B:122:ARG:HB3	1:B:136:HIS:HB2	1.67	0.74
1:B:189:TYR:O	3:B:302:HOH:O	2.06	0.73
1:A:181:GLN:N	3:A:307:HOH:O	2.21	0.73
1:A:181:GLN:OE1	3:A:302:HOH:O	2.08	0.70
1:A:194:ASP:O	3:A:304:HOH:O	2.09	0.69
1:B:11:TYR:N	3:B:304:HOH:O	2.18	0.69
1:A:21:ASP:OD1	3:A:305:HOH:O	2.10	0.69
1:A:100:THR:O	1:A:105:TYR:OH	2.10	0.68
1:B:85:TYR:HE2	1:B:133:THR:HG21	1.60	0.65
1:B:63:ARG:CZ	1:B:63:ARG:HA	2.27	0.65
1:A:114:TYR:O	3:A:306:HOH:O	2.15	0.65
1:A:178:VAL:O	3:A:307:HOH:O	2.15	0.65
1:A:159:LYS:HE3	1:A:160:LYS:HB3	1.80	0.64
1:A:119:ARG:HA	1:A:124:LYS:HG2	1.80	0.62
1:A:107:ARG:NH2	3:A:305:HOH:O	2.33	0.61
1:A:149:LEU:HB3	3:A:332:HOH:O	2.00	0.61
1:B:66:ASN:H	1:B:70:ILE:HD11	1.66	0.60
1:B:49:LYS:HA	1:B:98:LEU:HD11	1.83	0.60
1:A:47:VAL:HG11	1:A:157:LEU:HA	1.84	0.59
1:A:27:LEU:O	1:A:31:VAL:HG23	2.02	0.58
1:A:26:LEU:HB3	3:B:309:HOH:O	2.04	0.58
1:A:113:GLN:HG3	1:A:181:GLN:HG2	1.86	0.58
1:A:114:TYR:N	3:A:301:HOH:O	2.37	0.57
1:B:60:VAL:HB	2:B:201:WUF:C04	2.36	0.56
1:A:110:LEU:HD12	1:A:184:ILE:HD13	1.88	0.56
1:B:11:TYR:CZ	1:B:103:ARG:HG2	2.40	0.56
1:B:114:TYR:HA	3:B:324:HOH:O	2.05	0.56
1:A:88:ARG:HH21	1:A:129:ASP:HA	1.71	0.56
1:A:180:LEU:O	1:A:184:ILE:HG13	2.06	0.55
1:A:82:ILE:HG22	1:A:84:GLU:OE1	2.04	0.55
1:B:59:LEU:HD13	1:B:149:LEU:HD22	1.87	0.55
1:A:49:LYS:HA	1:A:98:LEU:HD21	1.88	0.55
1:A:84:GLU:H	1:A:84:GLU:CD	2.14	0.55
1:A:194:ASP:HA	1:B:136:HIS:CE1	2.42	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:60:VAL:HA	1:A:66:ASN:HD21	1.72	0.55
1:B:181:GLN:HG3	3:B:306:HOH:O	2.06	0.55
1:A:122:ARG:HB3	1:A:136:HIS:HB2	1.89	0.54
1:A:177:SER:OG	1:A:181:GLN:NE2	2.40	0.54
1:A:47:VAL:HG22	1:A:159:LYS:HE2	1.90	0.54
1:B:59:LEU:HD12	1:B:93:PHE:CE1	2.43	0.53
1:B:109:TRP:O	1:B:113:GLN:HG2	2.08	0.53
1:A:44:ALA:HB2	1:A:170:MET:HB3	1.91	0.52
1:B:48:LEU:HD22	1:B:55:ILE:HD11	1.91	0.52
1:B:185:TRP:HA	3:B:324:HOH:O	2.09	0.52
1:B:56:LEU:HD13	1:B:90:ARG:HG2	1.90	0.52
1:B:145:GLN:HG3	1:B:149:LEU:HG	1.92	0.52
1:A:69:LEU:HB3	2:A:201:WUF:CBC	2.40	0.51
1:A:15:GLU:OE1	1:A:15:GLU:N	2.44	0.51
1:A:185:TRP:O	1:A:188:PRO:HD2	2.12	0.50
1:A:140:ARG:HG3	1:B:27:LEU:HD21	1.94	0.50
1:A:193:ASN:N	3:A:313:HOH:O	2.42	0.50
1:B:10:ASP:N	3:B:304:HOH:O	2.44	0.50
1:A:166:ASP:HB2	3:A:322:HOH:O	2.11	0.50
1:B:166:ASP:O	1:B:170:MET:HG3	2.12	0.49
1:A:55:ILE:HG12	1:A:156:PHE:CD1	2.47	0.49
1:A:184:ILE:HG21	3:A:301:HOH:O	2.10	0.49
1:B:68:HIS:NE2	3:B:309:HOH:O	2.35	0.49
1:B:63:ARG:HA	1:B:63:ARG:NE	2.28	0.48
1:B:115:GLU:O	1:B:119:ARG:HG3	2.13	0.48
1:B:29:LYS:HA	1:B:29:LYS:HE2	1.94	0.48
1:A:60:VAL:HG23	1:A:86:LEU:HD23	1.95	0.48
3:A:304:HOH:O	1:B:189:TYR:HD2	1.96	0.47
1:A:178:VAL:C	3:A:307:HOH:O	2.56	0.47
1:A:94:GLY:HA2	1:A:97:ILE:HD12	1.96	0.47
1:A:124:LYS:O	1:A:124:LYS:HG3	2.15	0.47
1:B:47:VAL:HA	1:B:160:LYS:HG3	1.96	0.47
1:A:115:GLU:O	1:A:119:ARG:HG3	2.14	0.47
1:A:121:HIS:NE2	1:B:194:ASP:OD2	2.41	0.47
1:A:27:LEU:HD21	1:B:140:ARG:HG3	1.96	0.47
1:B:183:ALA:O	1:B:186:SER:OG	2.33	0.46
1:A:19:ILE:O	1:A:107:ARG:NH2	2.48	0.46
1:A:147:TYR:CZ	1:B:176:LYS:HE2	2.51	0.46
1:B:145:GLN:HB3	2:B:201:WUF:CAB	2.46	0.46
1:B:37:ASP:OD1	1:B:176:LYS:HE3	2.16	0.46
1:B:76:PRO:HD3	1:B:133:THR:C	2.41	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:22:LEU:O	1:A:26:LEU:HD12	2.16	0.46
1:A:66:ASN:HB2	1:A:70:ILE:HD12	1.97	0.45
1:B:43:LYS:O	1:B:47:VAL:HG23	2.17	0.45
1:B:66:ASN:O	1:B:69:LEU:HB2	2.17	0.45
1:A:62:TRP:CD1	1:A:83:LYS:HE3	2.52	0.45
1:A:166:ASP:O	1:A:170:MET:HG3	2.17	0.45
1:B:191:LYS:N	3:B:302:HOH:O	2.50	0.45
1:A:154:LYS:HB2	1:A:171:TYR:CE1	2.52	0.44
1:B:82:ILE:HD13	1:B:131:VAL:HG11	2.00	0.44
1:A:153:ILE:HG13	1:A:174:TRP:CE2	2.53	0.44
1:A:74:SER:HA	1:A:82:ILE:HG13	1.99	0.44
1:B:41:LEU:HG	1:B:173:ALA:HB1	2.00	0.44
1:A:59:LEU:HD23	1:A:93:PHE:CE1	2.53	0.44
1:B:56:LEU:HD11	1:B:94:GLY:HA3	2.00	0.44
1:A:174:TRP:HB3	3:A:320:HOH:O	2.18	0.43
1:B:112:TYR:O	1:B:116:VAL:HG23	2.18	0.43
1:A:114:TYR:HA	3:A:319:HOH:O	2.17	0.43
1:B:66:ASN:HB3	1:B:69:LEU:HB2	2.00	0.43
1:B:67:GLU:HB3	1:B:68:HIS:H	1.61	0.43
1:A:119:ARG:HG2	1:A:124:LYS:HD3	2.00	0.43
2:A:201:WUF:CHD	2:A:201:WUF:C01	2.95	0.43
1:B:47:VAL:O	1:B:51:GLN:NE2	2.51	0.43
1:A:78:THR:OG1	1:A:80:GLU:HB2	2.19	0.43
1:B:139:LEU:HG	1:B:143:ILE:HD12	2.01	0.43
1:A:164:PRO:O	1:A:168:GLU:HG2	2.18	0.43
2:B:201:WUF:CHD	2:B:201:WUF:C01	2.92	0.43
1:A:41:LEU:HG	1:A:173:ALA:HB1	2.00	0.43
1:A:60:VAL:HA	1:A:66:ASN:ND2	2.33	0.43
1:A:82:ILE:HB	1:A:85:TYR:HB3	2.01	0.43
1:A:112:TYR:O	1:A:116:VAL:HG23	2.19	0.42
1:A:113:GLN:HB2	3:A:301:HOH:O	2.19	0.42
1:A:187:HIS:HB3	1:A:188:PRO:HD3	2.01	0.42
1:A:52:VAL:HA	1:A:55:ILE:HD12	2.01	0.42
1:A:24:PHE:CD1	1:A:110:LEU:HD13	2.54	0.42
1:A:157:LEU:H	1:A:157:LEU:HG	1.71	0.42
1:A:154:LYS:HB2	1:A:171:TYR:CD1	2.55	0.42
1:B:185:TRP:HH2	2:B:201:WUF:C1B	2.33	0.42
1:A:136:HIS:CE1	1:A:138:PRO:HA	2.55	0.42
1:B:40:TYR:CE1	1:B:166:ASP:HA	2.55	0.42
1:A:159:LYS:O	1:A:160:LYS:C	2.63	0.42
1:A:76:PRO:HD3	1:A:133:THR:C	2.45	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:120:HIS:HA	1:A:137:ILE:HD12	2.02	0.41
1:B:11:TYR:OH	1:B:103:ARG:HG2	2.20	0.41
3:A:304:HOH:O	1:B:139:LEU:HB2	2.19	0.41
1:B:123:SER:C	1:B:124:LYS:HG2	2.45	0.41
1:A:26:LEU:O	1:A:30:THR:OG1	2.39	0.41
1:B:124:LYS:O	3:B:305:HOH:O	2.22	0.41
1:A:88:ARG:HB2	1:A:129:ASP:OD1	2.21	0.41
1:B:82:ILE:HD13	1:B:131:VAL:CG1	2.52	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:311:HOH:O	3:A:331:HOH:O[2_545]	2.05	0.15
1:A:115:GLU:OE2	1:A:132:ARG:NH1[2_555]	2.08	0.12

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all EM entries.

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	183/195 (94%)	173 (94%)	5 (3%)	5 (3%)	4	6
1	B	183/195 (94%)	174 (95%)	9 (5%)	0	100	100
All	All	366/390 (94%)	347 (95%)	14 (4%)	5 (1%)	11	17

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	66	ASN
1	A	65	SER
1	A	67	GLU
1	A	64	ALA

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Mol	Chain	Res	Type
1	A	192	GLU

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 PDB entries followed by that with respect to all EM entries.

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	166/173 (96%)	161 (97%)	5 (3%)	36	64
1	B	166/173 (96%)	158 (95%)	8 (5%)	23	46
All	All	332/346 (96%)	319 (96%)	13 (4%)	30	55

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

Mol	Chain	Res	Type
1	A	31	VAL
1	A	84	GLU
1	A	108	GLU
1	A	157	LEU
1	A	186	SER
1	B	21	ASP
1	B	29	LYS
1	B	67	GLU
1	B	86	LEU
1	B	108	GLU
1	B	123	SER
1	B	153	ILE
1	B	159	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	113	GLN
1	A	181	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

2 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	WUF	A	201	1	57,58,58	1.52	10 (17%)	73,95,95	2.13	18 (24%)
2	WUF	B	201	1	57,58,58	1.53	10 (17%)	73,95,95	2.22	21 (28%)

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	WUF	A	201	1	-	6/14/68/68	0/1/9/9
2	WUF	B	201	1	-	6/14/68/68	0/1/9/9

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	201	WUF	C1D-ND	3.74	1.44	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	201	WUF	C1D-ND	3.54	1.43	1.38
2	B	201	WUF	C03-C02	-3.53	1.32	1.38
2	A	201	WUF	C03-C02	-3.37	1.32	1.38
2	A	201	WUF	CAB-C3B	3.19	1.55	1.47
2	B	201	WUF	CAB-C3B	3.14	1.55	1.47
2	A	201	WUF	C01-C02	3.01	1.52	1.48
2	A	201	WUF	C4A-NA	2.98	1.43	1.38
2	B	201	WUF	C4A-NA	2.83	1.43	1.38
2	B	201	WUF	C01-C02	2.82	1.52	1.48
2	B	201	WUF	C1A-NA	2.73	1.42	1.38
2	A	201	WUF	C4D-ND	2.57	1.42	1.38
2	B	201	WUF	C4D-ND	2.57	1.42	1.38
2	A	201	WUF	C1A-NA	2.53	1.42	1.38
2	B	201	WUF	CAC-C3C	2.53	1.54	1.47
2	A	201	WUF	CAC-C3C	2.46	1.54	1.47
2	A	201	WUF	FE-NB	2.32	2.14	2.01
2	B	201	WUF	FE-NB	2.26	2.13	2.01
2	B	201	WUF	CMB-C2B	2.17	1.55	1.50
2	A	201	WUF	CMB-C2B	2.14	1.55	1.50

All (39) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	201	WUF	C4A-NA-C1A	-7.43	97.94	105.84
2	A	201	WUF	C4A-NA-C1A	-7.28	98.10	105.84
2	B	201	WUF	C3C-C4C-NC	-5.76	105.19	109.48
2	A	201	WUF	C3C-C4C-NC	-5.69	105.24	109.48
2	B	201	WUF	C1C-NC-C4C	5.55	111.75	105.84
2	A	201	WUF	C1C-NC-C4C	5.52	111.72	105.84
2	B	201	WUF	C4D-CHA-C1A	-5.48	114.38	126.02
2	A	201	WUF	C4D-CHA-C1A	-5.28	114.79	126.02
2	A	201	WUF	C01-C02-C07	-4.75	117.00	121.84
2	B	201	WUF	C01-C02-C07	-4.67	117.08	121.84
2	B	201	WUF	C4D-ND-C1D	-4.30	100.62	106.08
2	B	201	WUF	C2C-C1C-NC	-4.19	105.02	109.84
2	A	201	WUF	C4D-ND-C1D	-4.18	100.77	106.08
2	A	201	WUF	C2C-C1C-NC	-4.08	105.15	109.84
2	B	201	WUF	C3C-C2C-C1C	3.72	109.20	106.41
2	A	201	WUF	C3C-C2C-C1C	3.27	108.87	106.41
2	B	201	WUF	C4B-NB-C1B	3.16	110.08	106.08
2	B	201	WUF	CHD-C1D-ND	3.15	127.72	124.41
2	B	201	WUF	CHB-C4A-NA	3.14	128.26	124.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	201	WUF	C4B-NB-C1B	2.93	109.80	106.08
2	A	201	WUF	CHB-C4A-NA	2.89	127.95	124.37
2	B	201	WUF	C05-C06-C07	-2.73	116.87	120.24
2	A	201	WUF	CHD-C1D-ND	2.69	127.23	124.41
2	A	201	WUF	C05-C06-C07	-2.65	116.97	120.24
2	B	201	WUF	CHD-C4C-NC	2.62	127.25	124.44
2	A	201	WUF	CHD-C4C-NC	2.56	127.19	124.44
2	B	201	WUF	C3D-C4D-ND	2.38	112.62	110.33
2	A	201	WUF	CBC-CAC-C3C	-2.36	115.75	127.53
2	B	201	WUF	CHA-C1A-NA	-2.33	121.92	124.44
2	B	201	WUF	CAA-C2A-C1A	2.31	128.72	124.70
2	B	201	WUF	CHB-C4A-C3A	-2.29	120.44	126.95
2	A	201	WUF	CHB-C4A-C3A	-2.24	120.58	126.95
2	B	201	WUF	CBC-CAC-C3C	-2.21	116.50	127.53
2	A	201	WUF	C3D-C4D-ND	2.19	112.44	110.33
2	B	201	WUF	C4A-C3A-C2A	2.18	109.27	106.98
2	A	201	WUF	CHD-C1D-C2D	-2.13	121.05	125.49
2	B	201	WUF	CHD-C1D-C2D	-2.13	121.06	125.49
2	A	201	WUF	C2A-C1A-NA	2.05	112.32	110.34
2	B	201	WUF	C1A-C2A-C3A	-2.03	103.93	106.89

There are no chirality outliers.

All (12) torsion outliers are listed below:

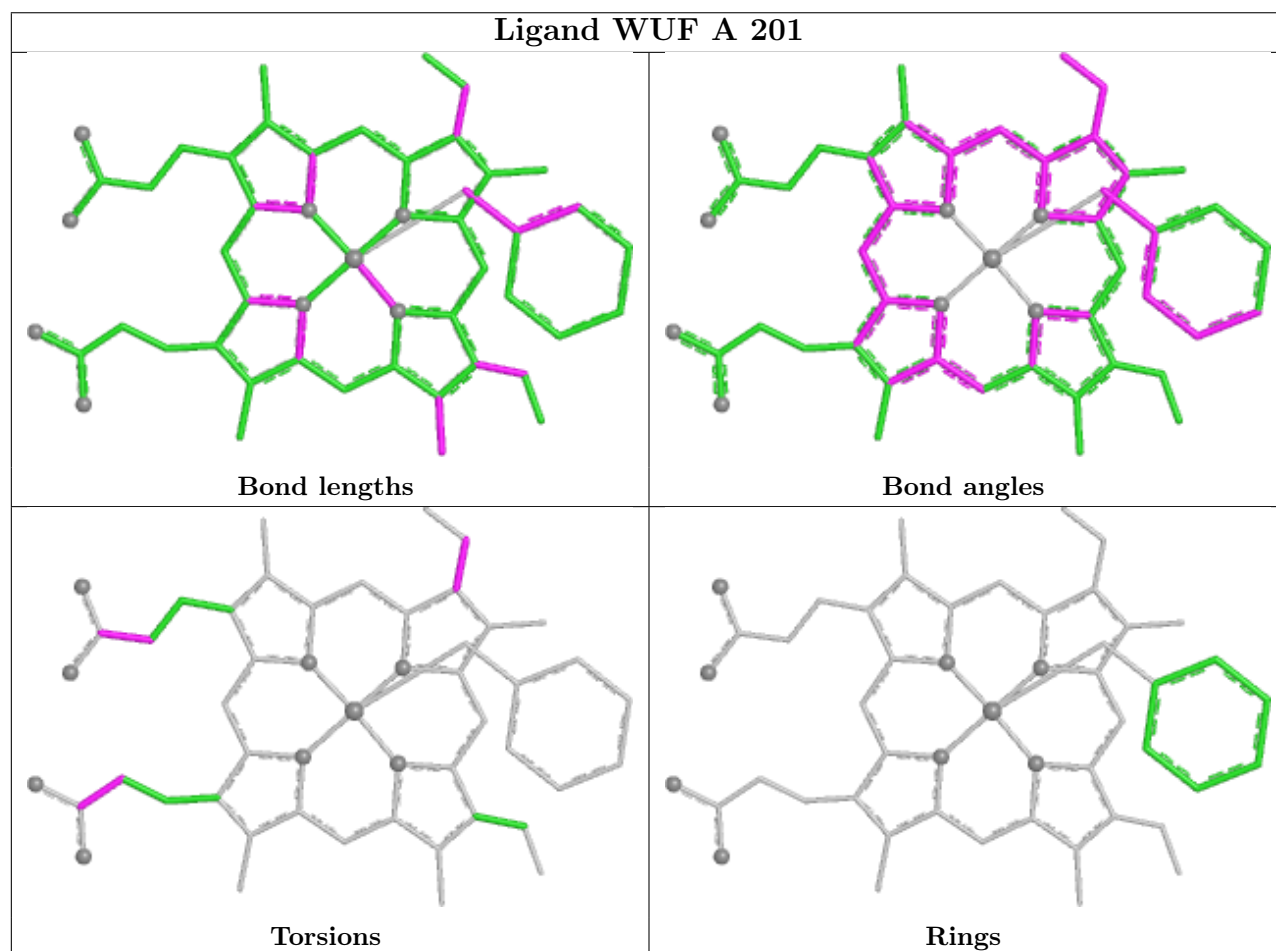
Mol	Chain	Res	Type	Atoms
2	A	201	WUF	C2C-C3C-CAC-CBC
2	B	201	WUF	C2C-C3C-CAC-CBC
2	A	201	WUF	C4C-C3C-CAC-CBC
2	B	201	WUF	C4C-C3C-CAC-CBC
2	A	201	WUF	CAD-CBD-CGD-O1D
2	B	201	WUF	CAD-CBD-CGD-O1D
2	A	201	WUF	CAA-CBA-CGA-O1A
2	B	201	WUF	CAA-CBA-CGA-O1A
2	A	201	WUF	CAA-CBA-CGA-O2A
2	B	201	WUF	CAA-CBA-CGA-O2A
2	A	201	WUF	CAD-CBD-CGD-O2D
2	B	201	WUF	CAD-CBD-CGD-O2D

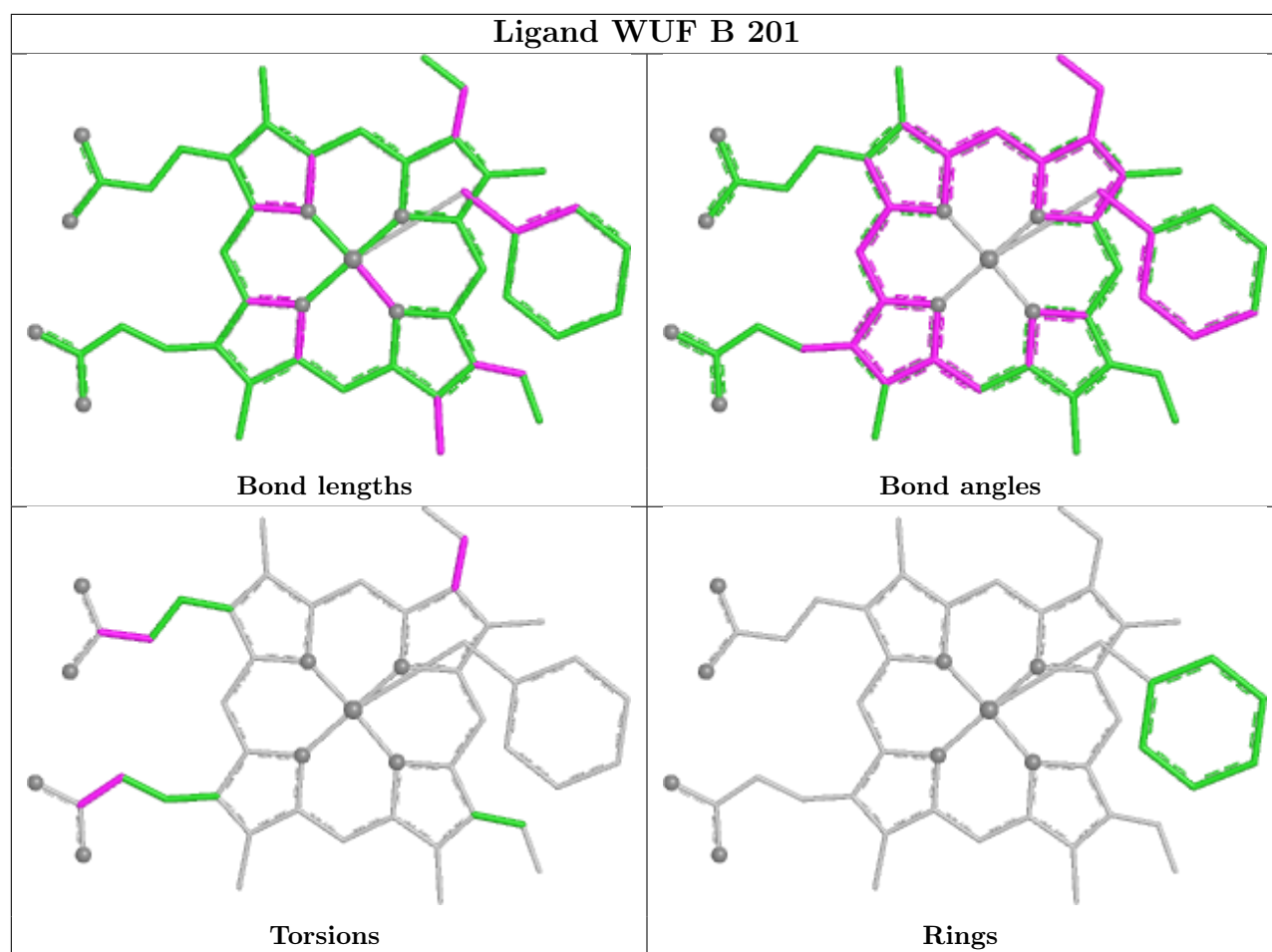
There are no ring outliers.

2 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	201	WUF	2	0
2	B	201	WUF	4	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.





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.