



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 4, 2026 – 10:45 AM EDT

PDB ID : 2CMM / pdb_00002cmm
Title : STRUCTURAL ANALYSIS OF THE MYOGLOBIN RECONSTITUTED
WITH IRON PORPHINE
Authors : Sato, T.; Tanaka, N.; Moriyama, H.; Igarashi, N.; Neya, S.; Funasaki, N.;
Iizuka, T.; Shiro, Y.
Deposited on : 1993-12-24
Resolution : 1.80 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)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.50

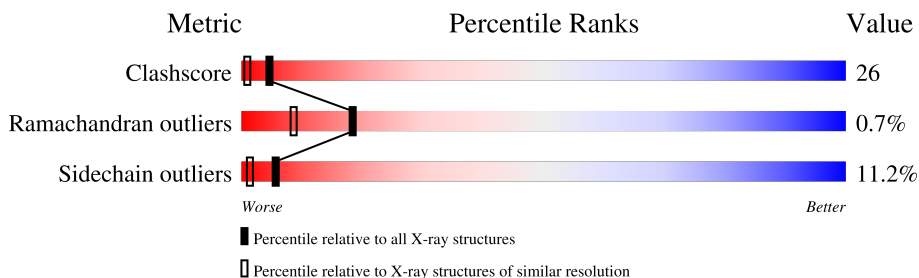
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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))
Clashscore	190562	8479 (1.80-1.80)
Ramachandran outliers	187476	8391 (1.80-1.80)
Sidechain outliers	187428	8390 (1.80-1.80)

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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	153	

2 Entry composition [i](#)

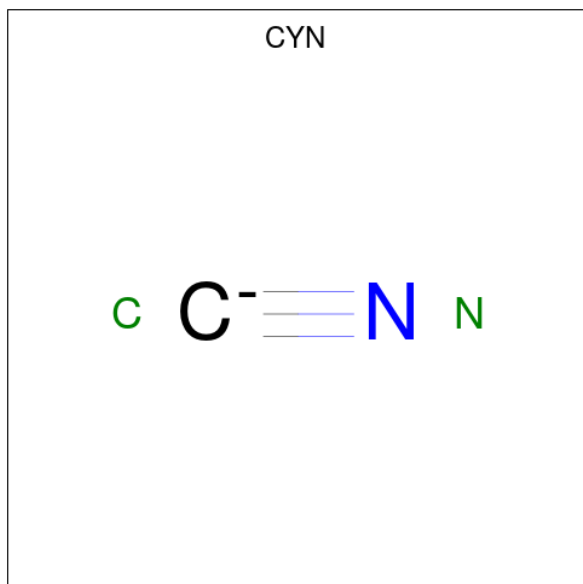
There are 4 unique types of molecules in this entry. The entry contains 1303 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 MYOGLOBIN.

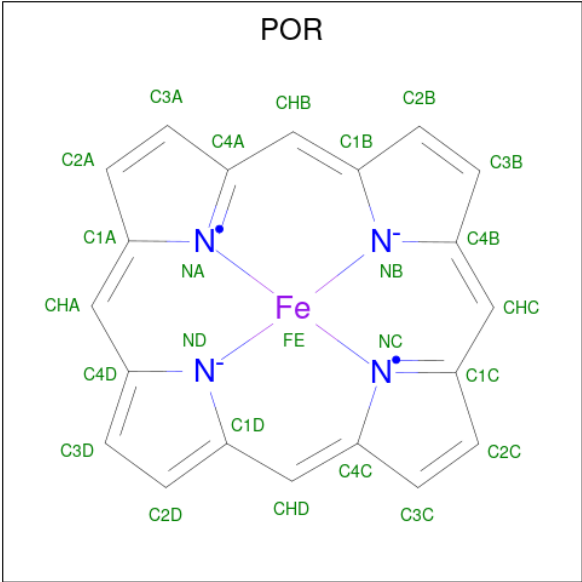
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	153	Total	C	N	O	S	0	0	0
			1217	783	216	216	2			

- Molecule 2 is CYANIDE ION (CCD ID: CYN) (formula: CN).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	N	0	0
			2	1	1		

- Molecule 3 is PORPHYRIN FE(III) (CCD ID: POR) (formula: C₂₀H₁₂FeN₄).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	Fe	N	0	0
			25	20	1	4		

- Molecule 4 is water.

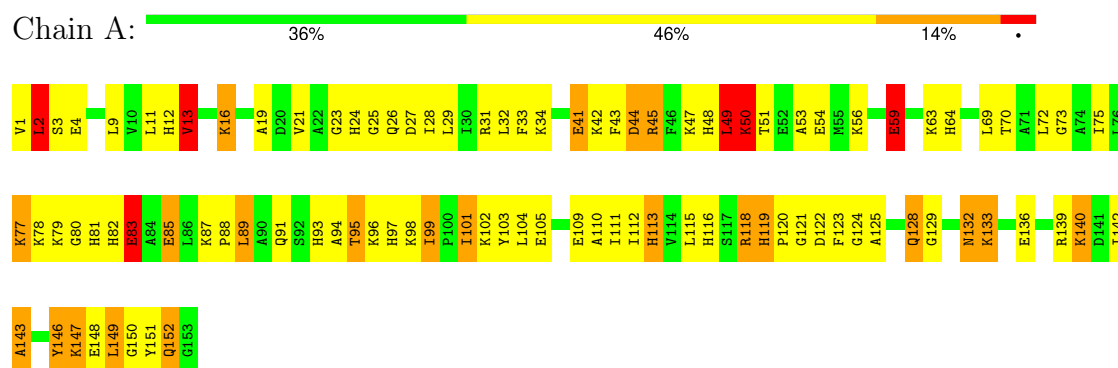
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	59	Total	O	0	0
			59	59		

3 Residue-property plots

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.

Note EDS was not executed.

• Molecule 1: MYOGLOBIN



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	57.62Å 76.23Å 33.26Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	(Not available) – 1.80	Depositor
% Data completeness (in resolution range)	(Not available) ((Not available)-1.80)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	PROFFT, X-PLOR	Depositor
R, R_{free}	0.187 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	1303	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: CYN, POR

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	1.78	15/1245 (1.2%)	2.62	99/1671 (5.9%)

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	26	GLN	N-CA	9.26	1.57	1.46
1	A	111	ILE	CA-CB	8.16	1.64	1.54
1	A	99	ILE	CA-CB	6.78	1.62	1.54
1	A	28	ILE	CA-CB	6.23	1.61	1.54
1	A	95	THR	CA-CB	6.22	1.63	1.53
1	A	110	ALA	C-N	5.78	1.41	1.33
1	A	129	GLY	CA-C	5.76	1.58	1.52
1	A	81	HIS	CA-CB	5.59	1.61	1.53
1	A	97	HIS	C-O	5.54	1.31	1.24
1	A	33	PHE	C-N	-5.43	1.26	1.33
1	A	115	LEU	C-N	-5.37	1.27	1.33
1	A	32	LEU	CA-CB	5.37	1.61	1.53
1	A	70	THR	N-CA	5.21	1.52	1.46
1	A	112	ILE	CB-CG1	5.20	1.63	1.53
1	A	125	ALA	CA-C	5.01	1.59	1.52

All (99) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	41	GLU	CA-CB-CG	12.78	139.66	114.10
1	A	81	HIS	CA-CB-CG	-11.93	101.87	113.80
1	A	83	GLU	CB-CG-CD	11.27	131.76	112.60
1	A	85	GLU	CB-CG-CD	11.16	131.57	112.60
1	A	139	ARG	NE-CZ-NH1	10.37	131.87	121.50
1	A	128	GLN	OE1-CD-NE2	10.08	132.68	122.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	119	HIS	CA-CB-CG	-9.80	104.00	113.80
1	A	96	LYS	CA-CB-CG	9.76	133.62	114.10
1	A	2	LEU	O-C-N	9.73	134.57	123.19
1	A	79	LYS	CB-CG-CD	9.58	133.34	111.30
1	A	113	HIS	CA-CB-CG	-9.49	104.31	113.80
1	A	133	LYS	CA-C-O	-9.15	110.85	120.55
1	A	125	ALA	O-C-N	8.79	131.12	122.07
1	A	96	LYS	N-CA-CB	8.60	123.52	110.19
1	A	116	HIS	CA-CB-CG	-8.14	105.66	113.80
1	A	132	ASN	CA-CB-CG	-7.92	104.68	112.60
1	A	125	ALA	CA-C-O	-7.75	112.68	120.82
1	A	27	ASP	O-C-N	-7.53	114.26	122.09
1	A	118	ARG	N-CA-C	7.27	120.63	111.69
1	A	23	GLY	CA-C-O	7.26	128.35	120.66
1	A	93	HIS	CA-CB-CG	-7.21	106.59	113.80
1	A	119	HIS	O-C-N	7.21	127.39	121.05
1	A	143	ALA	CA-C-O	-7.17	112.94	120.55
1	A	128	GLN	CA-C-N	7.15	127.88	119.94
1	A	128	GLN	C-N-CA	7.15	127.88	119.94
1	A	64	HIS	CA-C-N	7.05	127.80	119.98
1	A	64	HIS	C-N-CA	7.05	127.80	119.98
1	A	139	ARG	CD-NE-CZ	7.02	134.23	124.40
1	A	72	LEU	CA-C-N	7.02	127.77	119.98
1	A	72	LEU	C-N-CA	7.02	127.77	119.98
1	A	27	ASP	CA-C-O	6.98	128.37	120.90
1	A	96	LYS	CA-C-O	-6.84	111.76	120.31
1	A	95	THR	O-C-N	6.81	129.45	122.09
1	A	29	LEU	CA-C-O	-6.81	113.20	120.42
1	A	44	ASP	CA-CB-CG	-6.78	105.82	112.60
1	A	25	GLY	O-C-N	6.78	128.70	122.19
1	A	95	THR	N-CA-C	6.73	118.41	111.14
1	A	1	VAL	O-C-N	6.72	133.76	123.00
1	A	33	PHE	CA-C-O	-6.71	112.26	119.97
1	A	59	GLU	O-C-N	6.66	131.26	122.33
1	A	95	THR	CA-C-O	-6.63	113.87	120.70
1	A	29	LEU	O-C-N	6.49	129.55	122.15
1	A	45	ARG	CD-NE-CZ	-6.48	115.33	124.40
1	A	118	ARG	CA-C-O	-6.46	113.07	120.24
1	A	50	LYS	CB-CA-C	-6.43	102.47	112.31
1	A	69	LEU	O-C-N	6.40	130.14	122.27
1	A	3	SER	CA-C-O	6.29	129.16	121.87
1	A	132	ASN	O-C-N	6.27	128.53	122.07

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1	VAL	CA-C-O	-6.24	110.20	120.80
1	A	43	PHE	CA-C-O	6.22	127.12	120.46
1	A	33	PHE	O-C-N	6.02	129.68	122.27
1	A	26	GLN	OE1-CD-NE2	6.01	128.61	122.60
1	A	24	HIS	CA-C-N	5.99	126.63	119.98
1	A	24	HIS	C-N-CA	5.99	126.63	119.98
1	A	147	LYS	CA-C-O	5.97	127.29	120.90
1	A	118	ARG	CA-C-N	5.90	128.30	122.28
1	A	118	ARG	C-N-CA	5.90	128.30	122.28
1	A	41	GLU	N-CA-CB	5.85	120.24	110.41
1	A	12	HIS	CA-CB-CG	-5.84	107.95	113.80
1	A	129	GLY	O-C-N	5.83	127.77	122.18
1	A	1	VAL	CB-CA-C	-5.82	100.34	111.40
1	A	149	LEU	N-CA-C	-5.74	106.27	113.28
1	A	69	LEU	CA-C-O	-5.74	113.37	119.97
1	A	13	VAL	O-C-N	5.67	127.67	121.83
1	A	105	GLU	N-CA-C	-5.60	105.25	111.36
1	A	119	HIS	CB-CA-C	-5.60	103.95	111.11
1	A	136	GLU	OE1-CD-OE2	5.58	136.28	122.90
1	A	146	TYR	O-C-N	5.54	127.78	122.07
1	A	4	GLU	CB-CG-CD	5.53	122.00	112.60
1	A	3	SER	O-C-N	-5.53	116.49	122.79
1	A	2	LEU	N-CA-C	-5.50	100.74	109.76
1	A	31	ARG	CA-C-O	5.50	126.79	120.90
1	A	49	LEU	CA-C-O	-5.48	114.34	120.54
1	A	21	VAL	O-C-N	5.47	127.47	121.83
1	A	79	LYS	CA-CB-CG	-5.47	103.16	114.10
1	A	80	GLY	CA-C-O	-5.47	113.58	119.27
1	A	28	ILE	CA-CB-CG1	-5.41	101.21	110.40
1	A	49	LEU	CB-CA-C	5.38	118.63	109.80
1	A	150	GLY	N-CA-C	-5.38	107.84	115.43
1	A	4	GLU	CA-C-N	5.38	125.95	119.98
1	A	4	GLU	C-N-CA	5.38	125.95	119.98
1	A	128	GLN	N-CA-CB	5.33	117.96	110.12
1	A	121	GLY	CA-C-N	5.32	131.61	122.09
1	A	121	GLY	C-N-CA	5.32	131.61	122.09
1	A	101	ILE	CB-CG1-CD1	-5.31	102.65	113.80
1	A	70	THR	CA-CB-OG1	-5.30	101.65	109.60
1	A	149	LEU	CA-C-O	-5.30	112.87	119.18
1	A	16	LYS	CB-CG-CD	5.23	123.33	111.30
1	A	132	ASN	OD1-CG-ND2	5.21	127.81	122.60
1	A	2	LEU	CA-C-O	-5.21	115.01	120.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	59	GLU	CB-CG-CD	-5.21	103.74	112.60
1	A	148	GLU	CG-CD-OE2	-5.18	106.47	118.40
1	A	94	ALA	N-CA-CB	5.18	117.45	110.10
1	A	93	HIS	CE1-NE2-CD2	5.17	114.17	109.00
1	A	93	HIS	ND1-CE1-NE2	-5.17	103.23	108.40
1	A	59	GLU	CA-C-O	-5.17	113.69	119.79
1	A	139	ARG	NH1-CZ-NH2	-5.10	112.67	119.30
1	A	89	LEU	CA-C-N	5.08	127.36	120.65
1	A	89	LEU	C-N-CA	5.08	127.36	120.65

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	1217	0	1241	65	0
2	A	2	0	0	0	0
3	A	25	0	12	0	0
4	A	59	0	0	6	0
All	All	1303	0	1253	65	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 26.

All (65) 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:78:LYS:HE2	1:A:82:HIS:HA	1.04	1.02
1:A:123:PHE:CE2	1:A:128:GLN:HG2	1.99	0.97
1:A:152:GLN:O	4:A:251:HOH:O	1.82	0.95
1:A:78:LYS:CE	1:A:82:HIS:HA	1.97	0.95
1:A:51:THR:HG23	1:A:54:GLU:OE2	1.65	0.94
1:A:140:LYS:HZ3	1:A:140:LYS:HB2	1.42	0.83
1:A:78:LYS:NZ	1:A:85:GLU:H	1.79	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:78:LYS:HE2	1:A:82:HIS:CA	2.00	0.80
1:A:19:ALA:O	4:A:242:HOH:O	2.00	0.78
1:A:78:LYS:HZ3	1:A:85:GLU:H	1.34	0.74
1:A:42:LYS:HG2	1:A:103:TYR:HE2	1.50	0.74
1:A:16:LYS:HE2	1:A:122:ASP:OD2	1.87	0.73
1:A:42:LYS:CG	1:A:103:TYR:HE2	2.02	0.73
1:A:59:GLU:CD	1:A:63:LYS:HE2	2.16	0.70
1:A:95:THR:OG1	4:A:215:HOH:O	2.11	0.69
1:A:78:LYS:HD2	1:A:85:GLU:CD	2.19	0.67
1:A:140:LYS:HB2	1:A:140:LYS:NZ	2.09	0.66
1:A:152:GLN:NE2	1:A:152:GLN:HA	2.12	0.63
1:A:51:THR:OG1	1:A:53:ALA:HB3	1.99	0.62
1:A:143:ALA:O	1:A:147:LYS:HD3	1.99	0.61
1:A:87:LYS:HB3	1:A:88:PRO:HD3	1.82	0.61
1:A:78:LYS:NZ	1:A:85:GLU:HB2	2.16	0.60
1:A:78:LYS:NZ	1:A:85:GLU:CB	2.65	0.60
1:A:123:PHE:CD2	1:A:128:GLN:HG2	2.35	0.60
1:A:49:LEU:C	1:A:50:LYS:HD2	2.28	0.59
1:A:140:LYS:NZ	1:A:140:LYS:CB	2.66	0.59
1:A:75:ILE:HD11	1:A:89:LEU:HD23	1.84	0.58
1:A:59:GLU:OE1	1:A:63:LYS:HE2	2.03	0.58
1:A:83:GLU:H	1:A:83:GLU:CD	2.12	0.57
1:A:113:HIS:CD2	1:A:113:HIS:C	2.81	0.57
1:A:44:ASP:O	1:A:47:LYS:HG2	2.05	0.57
1:A:124:GLY:O	1:A:128:GLN:HG3	2.07	0.55
1:A:78:LYS:NZ	1:A:85:GLU:N	2.54	0.54
1:A:78:LYS:HZ3	1:A:85:GLU:HB2	1.72	0.54
1:A:50:LYS:HD2	1:A:50:LYS:N	2.23	0.54
1:A:42:LYS:HG3	1:A:103:TYR:CE2	2.43	0.54
1:A:42:LYS:CG	1:A:103:TYR:CE2	2.88	0.53
1:A:50:LYS:N	1:A:54:GLU:OE1	2.39	0.53
1:A:13:VAL:O	1:A:16:LYS:HB2	2.10	0.53
1:A:119:HIS:N	1:A:120:PRO:HD3	2.26	0.51
1:A:73:GLY:O	1:A:77:LYS:HD2	2.11	0.50
1:A:51:THR:O	1:A:54:GLU:HB2	2.11	0.50
1:A:152:GLN:NE2	1:A:152:GLN:CA	2.77	0.48
1:A:51:THR:HG1	1:A:53:ALA:HB3	1.78	0.47
1:A:104:LEU:HD21	1:A:142:ILE:HG21	1.96	0.47
1:A:45:ARG:HH11	1:A:45:ARG:HD3	1.49	0.47
1:A:132:ASN:HB2	4:A:256:HOH:O	2.15	0.46
1:A:118:ARG:C	1:A:120:PRO:HD3	2.41	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:48:HIS:H	1:A:48:HIS:CD2	2.32	0.46
1:A:78:LYS:HD2	1:A:85:GLU:CG	2.45	0.46
1:A:2:LEU:CD2	1:A:133:LYS:HB3	2.46	0.46
1:A:59:GLU:O	1:A:63:LYS:HG3	2.15	0.46
1:A:140:LYS:HZ3	1:A:140:LYS:CB	2.18	0.45
1:A:34:LYS:HD3	1:A:34:LYS:HA	1.73	0.44
1:A:152:GLN:C	4:A:251:HOH:O	2.52	0.44
1:A:101:ILE:HD12	1:A:101:ILE:HG23	1.69	0.43
1:A:91:GLN:HG2	1:A:149:LEU:HD11	2.01	0.42
1:A:59:GLU:OE2	1:A:63:LYS:NZ	2.47	0.42
1:A:51:THR:H	1:A:54:GLU:CD	2.28	0.42
1:A:99:ILE:HG23	1:A:103:TYR:HB2	2.02	0.42
1:A:146:TYR:CD2	1:A:151:TYR:HD2	2.38	0.42
1:A:101:ILE:HD13	1:A:101:ILE:HA	1.89	0.41
1:A:109:GLU:CD	4:A:269:HOH:O	2.64	0.41
1:A:119:HIS:HB3	1:A:122:ASP:HB2	2.03	0.40
1:A:2:LEU:HD22	1:A:133:LYS:HB3	2.03	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	151/153 (99%)	145 (96%)	5 (3%)	1 (1%)	18 8

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	152	GLN

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all 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	125/125 (100%)	111 (89%)	14 (11%)	6 1

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	LEU
1	A	9	LEU
1	A	11	LEU
1	A	13	VAL
1	A	41	GLU
1	A	49	LEU
1	A	50	LYS
1	A	56	LYS
1	A	59	GLU
1	A	77	LYS
1	A	83	GLU
1	A	98	LYS
1	A	102	LYS
1	A	140	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	91	GLN
1	A	113	HIS
1	A	128	GLN
1	A	132	ASN
1	A	152	GLN

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](#)

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	CYN	A	154	-	1,1,1	1.23	0	-		
3	POR	A	155	1	32,32,32	1.18	4 (12%)	30,54,54	1.78	5 (16%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	155	POR	FE-NC	2.54	2.02	1.94
3	A	155	POR	C3D-C4D	2.08	1.44	1.40
3	A	155	POR	C1C-NC	2.01	1.41	1.38
3	A	155	POR	FE-NA	2.00	2.01	1.94

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	155	POR	C3A-C4A-NA	-7.26	103.96	110.01
3	A	155	POR	CHB-C4A-NA	-3.01	121.17	124.78
3	A	155	POR	C3D-C4D-ND	-2.79	107.35	109.71
3	A	155	POR	C2A-C3A-C4A	2.14	109.94	107.24
3	A	155	POR	C2D-C3D-C4D	2.02	109.05	106.81

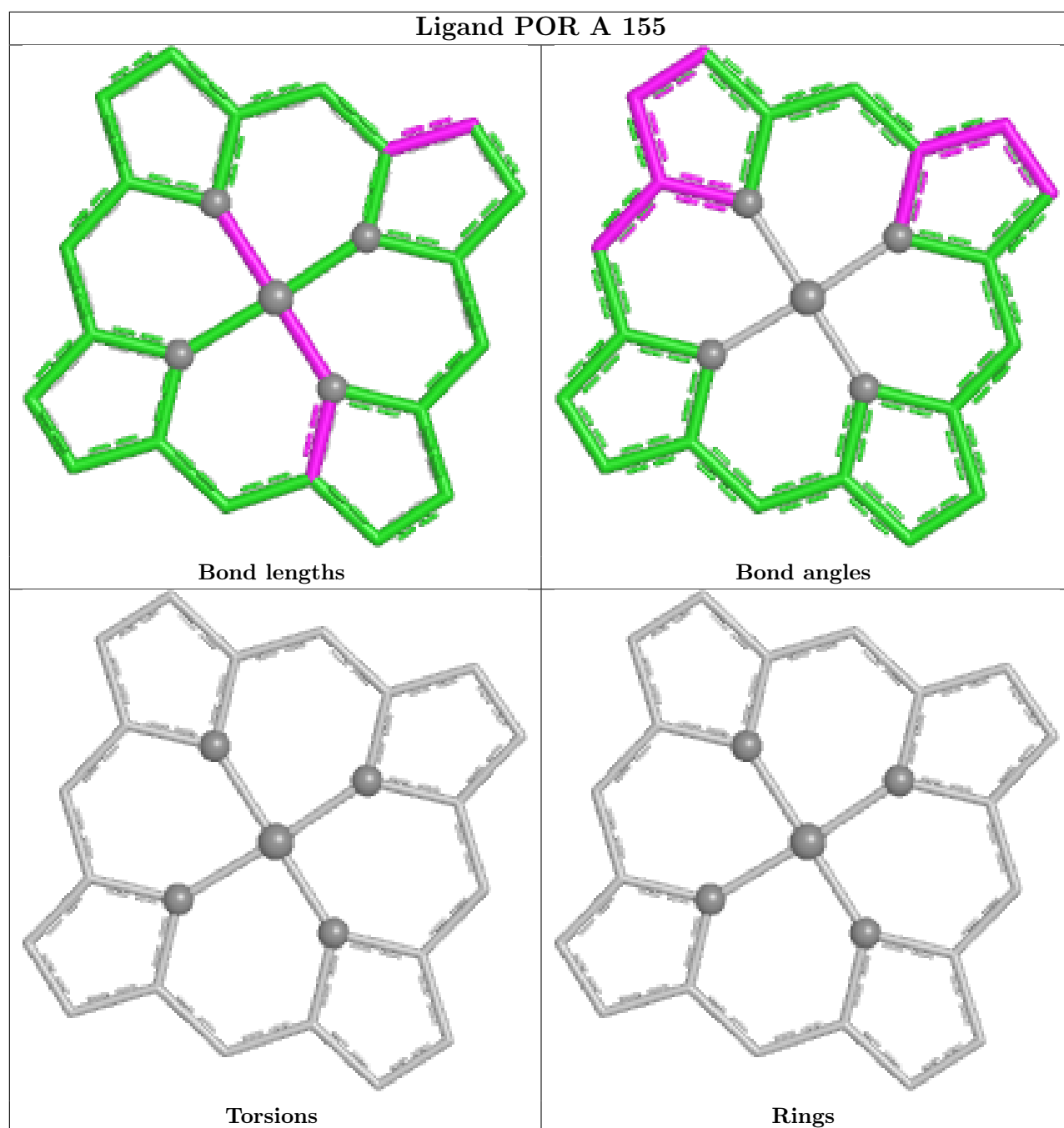
There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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

6 Fit of model and data

6.1 Protein, DNA and RNA chains

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.