



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 4, 2026 – 08:51 PM EDT

PDB ID : 2D2M / pdb_00002d2m
Title : Structure of an extracellular giant hemoglobin of the gutless beard worm *Oligobrachia mashikoi*
Authors : Numoto, N.; Nakagawa, T.; Kita, A.; Sasayama, Y.; Fukumori, Y.; Miki, K.
Deposited on : 2005-09-12
Resolution : 2.85 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

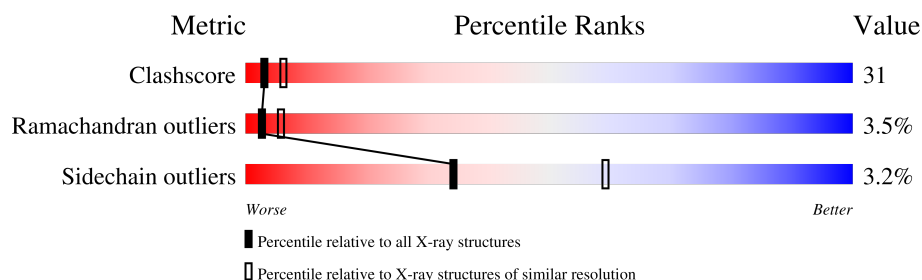
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.85 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	1446 (2.88-2.84)
Ramachandran outliers	187476	1406 (2.88-2.84)
Sidechain outliers	187428	1407 (2.88-2.84)

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	140	41% 53% 6%
2	B	142	49% 47% .
3	C	147	46% 47% 6% ..
4	D	145	50% 47% ..

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 4457 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 Giant hemoglobin, A1(b) globin chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	140	Total	C	N	O	S	0	0	0
			1065	656	196	206	7			

- Molecule 2 is a protein called Giant hemoglobin, A2(a5) globin chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	142	Total	C	N	O	S	0	0	0
			1075	662	198	209	6			

- Molecule 3 is a protein called Giant hemoglobin, B2(c) globin chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	146	Total	C	N	O	S	0	0	0
			1088	668	201	210	9			

- Molecule 4 is a protein called Giant hemoglobin, B1(d) globin chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	145	Total	C	N	O	S	0	0	0
			1036	642	178	209	7			

- Molecule 5 is PROTOPORPHYRIN IX CONTAINING FE (CCD ID: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
5	B	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
5	C	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
5	D	1	Total 43	C 34	Fe 1	N 4	O 4	0	0

- Molecule 6 is OXYGEN MOLECULE (CCD ID: OXY) (formula: O_2).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	1	Total O 2 2	0	0
6	B	1	Total O 2 2	0	0
6	C	1	Total O 2 2	0	0
6	D	1	Total O 2 2	0	0

- Molecule 7 is water.

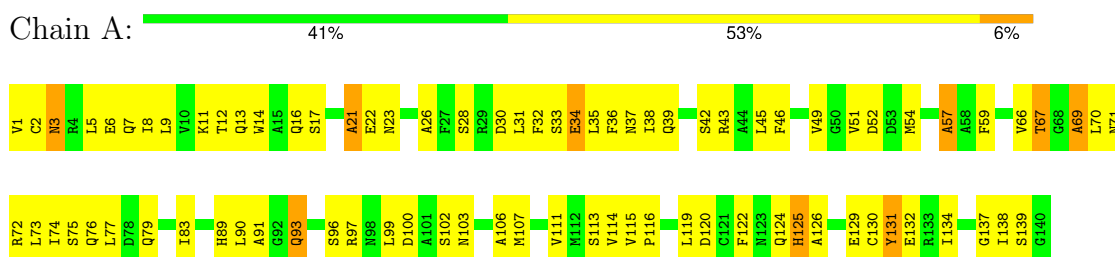
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	6	Total O 6 6	0	0
7	B	1	Total O 1 1	0	0
7	C	3	Total O 3 3	0	0
7	D	3	Total O 3 3	0	0

3 Residue-property plots

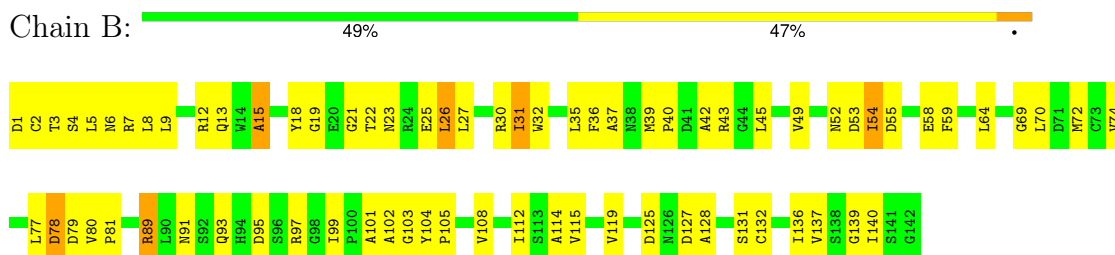
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded 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.

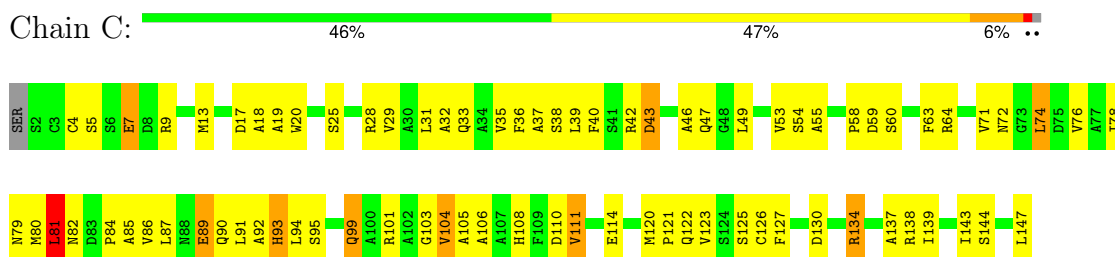
- Molecule 1: Giant hemoglobin, A1(b) globin chain



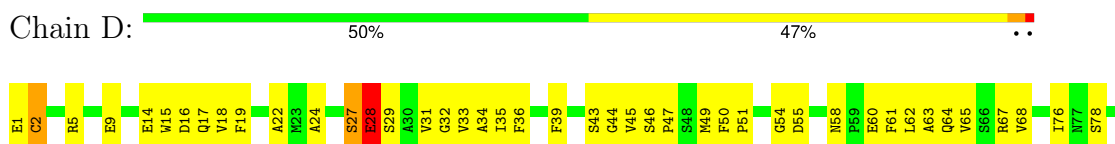
- Molecule 2: Giant hemoglobin, A2(a5) globin chain



- Molecule 3: Giant hemoglobin, B2(c) globin chain



- Molecule 4: Giant hemoglobin, B1(d) globin chain



T84		L89	S90		N93		R97		V102		A106	V107	T108	H109	L110	S111	Q112	A113	I114	S115	S116	V117	V118		V121	L122	P123	S124	A125	H126	I127	D128	A129	W130		M134		I137		G142		L145
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4 Data and refinement statistics

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

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants a, b, c, α , β , γ	111.50Å 111.50Å 276.83Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	50.00 – 2.85	Depositor
% Data completeness (in resolution range)	97.3 (50.00-2.85)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.236 , 0.277	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	4457	wwPDB-VP
Average B, all atoms (Å ²)	58.0	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: OXY, HEM

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	0.41	0/1083	0.90	8/1464 (0.5%)
2	B	0.43	0/1093	0.83	1/1481 (0.1%)
3	C	0.41	0/1110	0.89	3/1507 (0.2%)
4	D	0.40	0/1055	0.90	5/1436 (0.3%)
All	All	0.41	0/4341	0.88	17/5888 (0.3%)

There are no bond length outliers.

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	120	ASP	N-CA-C	-7.76	103.94	113.41
4	D	28	GLU	N-CA-C	-7.22	103.73	112.54
3	C	99	GLN	N-CA-C	-5.69	105.16	111.36
4	D	16	ASP	N-CA-C	-5.67	106.45	113.19
1	A	113	SER	N-CA-C	-5.57	106.85	113.97
2	B	15	ALA	N-CA-C	-5.39	105.09	110.97
1	A	3	ASN	N-CA-C	5.38	117.03	110.41
4	D	27	SER	N-CA-C	5.35	116.86	110.44
1	A	69	ALA	N-CA-C	-5.26	105.44	111.07
1	A	57	ALA	N-CA-C	-5.26	105.63	111.36
1	A	67	THR	N-CA-C	-5.25	105.98	112.38
1	A	131	TYR	N-CA-C	-5.19	105.62	111.28
3	C	17	ASP	N-CA-C	-5.11	105.62	111.14
1	A	93	GLN	N-CA-C	-5.08	107.03	113.18
4	D	18	VAL	N-CA-C	-5.02	104.81	111.09
4	D	44	GLY	N-CA-C	-5.00	109.06	115.42
3	C	111	VAL	N-CA-C	-5.00	105.67	110.72

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1065	0	1015	78	0
2	B	1075	0	1038	77	0
3	C	1088	0	1013	75	0
4	D	1036	0	987	58	0
5	A	43	0	30	4	0
5	B	43	0	30	4	0
5	C	43	0	30	0	0
5	D	43	0	30	2	0
6	A	2	0	0	0	0
6	B	2	0	0	0	0
6	C	2	0	0	0	0
6	D	2	0	0	0	0
7	A	6	0	0	0	0
7	B	1	0	0	0	0
7	C	3	0	0	0	0
7	D	3	0	0	0	0
All	All	4457	0	4173	271	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 31.

All (271) 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)
3:C:25:SER:HA	3:C:28:ARG:HD3	1.42	0.98
1:A:14:TRP:HE1	1:A:71:ASN:ND2	1.65	0.95
1:A:46:PHE:HB2	1:A:51:VAL:HG11	1.50	0.93
2:B:9:LEU:HD21	2:B:13:GLN:HE21	1.34	0.93
1:A:14:TRP:NE1	1:A:71:ASN:HD22	1.69	0.91
3:C:19:ALA:HB1	3:C:123:VAL:HG12	1.54	0.89
1:A:14:TRP:HE1	1:A:71:ASN:HD22	0.85	0.84
2:B:69:GLY:HA2	2:B:72:MET:HE2	1.58	0.83
3:C:49:LEU:HD11	3:C:108:HIS:CE1	2.16	0.80
1:A:72:ARG:NE	4:D:67:ARG:HG2	1.97	0.80
2:B:21:GLY:H	3:C:79:ASN:HD22	1.32	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:39:GLN:HE22	1:A:103:ASN:HA	1.49	0.76
1:A:138:ILE:HD12	5:A:200:HEM:HBB2	1.69	0.75
4:D:51:PRO:HD2	4:D:64:GLN:HG2	1.68	0.75
1:A:100:ASP:H	1:A:103:ASN:ND2	1.85	0.74
4:D:1:GLU:HG3	4:D:2:CYS:H	1.52	0.73
1:A:91:ALA:HB2	1:A:138:ILE:HA	1.71	0.73
4:D:61:PHE:O	4:D:65:VAL:HG23	1.90	0.72
4:D:110:LEU:HD12	5:D:200:HEM:HBB2	1.72	0.70
3:C:58:PRO:HA	3:C:63:PHE:CD1	2.28	0.69
2:B:4:SER:HA	2:B:7:ARG:HE	1.58	0.69
2:B:97:ARG:HG2	5:B:200:HEM:HMD3	1.74	0.68
3:C:4:CYS:SG	3:C:9:ARG:HG2	2.34	0.68
1:A:7:GLN:O	1:A:11:LYS:HG3	1.93	0.68
3:C:25:SER:HA	3:C:28:ARG:CD	2.23	0.68
2:B:26:LEU:HD11	2:B:30:ARG:HH21	1.59	0.67
1:A:46:PHE:HB2	1:A:51:VAL:CG1	2.22	0.67
3:C:32:ALA:HB2	3:C:71:VAL:CG2	2.24	0.67
2:B:35:LEU:HD11	2:B:39:MET:HE2	1.76	0.67
2:B:115:VAL:O	2:B:119:VAL:HG22	1.94	0.66
2:B:97:ARG:H	2:B:97:ARG:HD2	1.59	0.66
1:A:8:ILE:HD13	3:C:123:VAL:HG13	1.77	0.66
2:B:1:ASP:CG	2:B:2:CYS:H	2.04	0.66
2:B:89:ARG:O	2:B:93:GLN:HG3	1.96	0.65
3:C:91:LEU:HD21	3:C:139:ILE:HG23	1.79	0.65
4:D:35:ILE:HD12	4:D:114:ILE:HA	1.77	0.65
2:B:23:ASN:O	2:B:26:LEU:HB3	1.97	0.64
4:D:35:ILE:HD13	4:D:117:VAL:HG21	1.80	0.64
3:C:28:ARG:HB3	3:C:71:VAL:HG11	1.80	0.64
4:D:64:GLN:O	4:D:68:VAL:HG23	1.98	0.63
1:A:72:ARG:HH22	1:A:90:LEU:HD21	1.63	0.63
3:C:5:SER:O	3:C:9:ARG:HG3	1.99	0.63
3:C:76:VAL:O	3:C:80:MET:HG3	1.99	0.63
1:A:12:THR:O	1:A:16:GLN:HG3	1.98	0.62
2:B:99:ILE:HD12	5:B:200:HEM:HAC	1.80	0.61
3:C:90:GLN:HE21	3:C:90:GLN:HA	1.63	0.61
3:C:72:ASN:O	3:C:76:VAL:HG23	2.00	0.61
3:C:9:ARG:HD3	3:C:81:LEU:O	1.99	0.61
2:B:99:ILE:HG23	5:B:200:HEM:HAC	1.83	0.61
3:C:32:ALA:HB2	3:C:71:VAL:HG23	1.82	0.60
1:A:97:ARG:H	1:A:97:ARG:HD2	1.67	0.60
3:C:53:VAL:CG1	3:C:63:PHE:HA	2.30	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:74:LEU:HD22	3:C:78:ILE:HG13	1.84	0.59
3:C:39:LEU:HD12	3:C:111:VAL:HG12	1.86	0.58
2:B:4:SER:CB	2:B:7:ARG:HH21	2.17	0.58
1:A:107:MET:HB3	1:A:131:TYR:OH	2.03	0.57
2:B:58:GLU:HA	3:C:89:GLU:HG2	1.85	0.57
1:A:97:ARG:HG2	5:A:200:HEM:HMD1	1.86	0.57
4:D:110:LEU:HD13	4:D:110:LEU:O	2.04	0.57
2:B:80:VAL:HB	2:B:81:PRO:HD3	1.85	0.57
1:A:72:ARG:HE	4:D:67:ARG:HG2	1.70	0.56
2:B:39:MET:HE3	2:B:42:ALA:HB2	1.86	0.56
4:D:36:PHE:HZ	4:D:68:VAL:HG21	1.68	0.56
3:C:29:VAL:O	3:C:33:GLN:HG3	2.06	0.56
2:B:43:ARG:HB2	2:B:43:ARG:NH1	2.21	0.56
2:B:43:ARG:HB2	2:B:43:ARG:HH11	1.71	0.56
1:A:116:PRO:HB3	1:A:122:PHE:CD2	2.40	0.56
2:B:9:LEU:CD2	2:B:13:GLN:HE21	2.12	0.56
1:A:11:LYS:HE3	1:A:75:SER:HA	1.88	0.56
2:B:5:LEU:HD23	2:B:8:LEU:HD12	1.87	0.55
1:A:1:VAL:HG12	1:A:129:GLU:HG3	1.89	0.55
4:D:114:ILE:HB	4:D:134:MET:HE2	1.88	0.55
1:A:5:LEU:HD12	3:C:122:GLN:NE2	2.22	0.55
3:C:82:ASN:C	3:C:84:PRO:HD3	2.32	0.55
3:C:92:ALA:O	3:C:95:SER:HB3	2.06	0.55
4:D:115:SER:OG	4:D:134:MET:HG3	2.07	0.55
4:D:35:ILE:HD13	4:D:117:VAL:CG2	2.36	0.54
2:B:31:ILE:HG22	2:B:114:ALA:CB	2.37	0.54
2:B:39:MET:HE3	2:B:42:ALA:CB	2.38	0.54
2:B:42:ALA:O	2:B:45:LEU:HD23	2.08	0.54
1:A:138:ILE:O	1:A:138:ILE:HG22	2.08	0.54
2:B:9:LEU:HD23	2:B:9:LEU:C	2.32	0.54
1:A:74:ILE:HA	1:A:77:LEU:HG	1.89	0.54
1:A:125:HIS:O	1:A:126:ALA:C	2.51	0.54
1:A:54:MET:HA	1:A:59:PHE:CD2	2.43	0.54
4:D:108:THR:HA	4:D:111:SER:OG	2.06	0.54
1:A:72:ARG:HG3	1:A:73:LEU:HD12	1.89	0.54
4:D:19:PHE:HB3	4:D:24:ALA:HB2	1.90	0.54
1:A:45:LEU:HD21	1:A:99:LEU:HD23	1.89	0.53
2:B:3:THR:H	2:B:6:ASN:HD22	1.55	0.53
1:A:30:ASP:O	1:A:33:SER:HB3	2.09	0.53
2:B:26:LEU:HD13	2:B:26:LEU:C	2.33	0.53
2:B:70:LEU:O	2:B:74:VAL:HG23	2.09	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:21:GLY:H	3:C:79:ASN:ND2	2.04	0.53
1:A:6:GLU:OE1	1:A:129:GLU:HG2	2.08	0.53
1:A:43:ARG:HH11	1:A:43:ARG:HG2	1.74	0.53
2:B:36:PHE:HB3	2:B:43:ARG:HH12	1.74	0.53
2:B:69:GLY:HA2	2:B:72:MET:CE	2.33	0.53
3:C:103:GLY:O	3:C:104:VAL:C	2.51	0.53
4:D:19:PHE:CE1	4:D:28:GLU:HG3	2.44	0.52
3:C:53:VAL:O	3:C:55:ALA:N	2.34	0.52
1:A:28:SER:HA	1:A:67:THR:HG21	1.91	0.52
1:A:34:GLU:HA	1:A:34:GLU:OE1	2.10	0.52
3:C:38:SER:O	3:C:39:LEU:C	2.53	0.52
2:B:9:LEU:HD21	2:B:13:GLN:NE2	2.14	0.52
1:A:5:LEU:HA	1:A:8:ILE:CD1	2.40	0.52
2:B:3:THR:N	2:B:6:ASN:HD22	2.08	0.52
3:C:90:GLN:HA	3:C:90:GLN:NE2	2.25	0.51
1:A:2:CYS:HB2	1:A:130:CYS:HA	1.91	0.51
4:D:33:VAL:HG23	4:D:34:ALA:N	2.25	0.51
2:B:64:LEU:HD12	3:C:80:MET:HE2	1.92	0.51
1:A:51:VAL:HG12	1:A:59:PHE:CE1	2.46	0.51
2:B:108:VAL:CG2	2:B:137:VAL:HG21	2.41	0.51
3:C:31:LEU:O	3:C:35:VAL:HG23	2.11	0.51
1:A:17:SER:HB3	1:A:119:LEU:HD21	1.93	0.51
2:B:35:LEU:HD11	2:B:39:MET:CE	2.41	0.51
1:A:39:GLN:HE21	1:A:42:SER:HB3	1.75	0.50
2:B:43:ARG:HD3	2:B:52:ASN:HA	1.92	0.50
2:B:7:ARG:NH1	2:B:78:ASP:HA	2.25	0.50
1:A:91:ALA:HB2	1:A:137:GLY:O	2.11	0.50
4:D:1:GLU:HG3	4:D:2:CYS:N	2.23	0.50
1:A:96:SER:OG	1:A:97:ARG:HD2	2.12	0.50
3:C:94:LEU:HD13	3:C:143:ILE:HD11	1.93	0.50
2:B:52:ASN:N	2:B:52:ASN:HD22	2.09	0.49
2:B:91:ASN:HB2	2:B:139:GLY:O	2.11	0.49
2:B:5:LEU:HD22	4:D:121:VAL:CG2	2.42	0.49
1:A:36:PHE:CD1	1:A:43:ARG:HG3	2.48	0.49
1:A:97:ARG:HD2	1:A:97:ARG:N	2.26	0.49
4:D:61:PHE:O	4:D:64:GLN:HB3	2.12	0.49
3:C:9:ARG:O	3:C:13:MET:HG3	2.13	0.49
3:C:120:MET:HB2	3:C:121:PRO:HD3	1.94	0.49
3:C:137:ALA:O	3:C:138:ARG:C	2.56	0.49
4:D:102:VAL:HG22	5:D:200:HEM:HAC	1.94	0.49
1:A:38:ILE:HD11	1:A:106:ALA:O	2.12	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:69:ALA:O	1:A:73:LEU:HD13	2.13	0.48
3:C:25:SER:CA	3:C:28:ARG:HD3	2.30	0.48
2:B:22:THR:O	2:B:25:GLU:HB2	2.12	0.48
4:D:36:PHE:O	4:D:39:PHE:HB3	2.13	0.48
2:B:1:ASP:CG	2:B:2:CYS:N	2.71	0.48
1:A:32:PHE:HZ	1:A:66:VAL:HG21	1.78	0.48
1:A:35:LEU:HD13	1:A:107:MET:HA	1.96	0.48
3:C:19:ALA:CB	3:C:123:VAL:HG12	2.35	0.48
3:C:39:LEU:HD12	3:C:111:VAL:CG1	2.43	0.48
4:D:31:VAL:O	4:D:35:ILE:HG12	2.14	0.48
2:B:74:VAL:HA	2:B:77:LEU:HG	1.96	0.47
4:D:47:PRO:C	4:D:49:MET:H	2.22	0.47
2:B:101:ALA:C	2:B:103:GLY:H	2.22	0.47
4:D:107:VAL:HG21	4:D:142:GLY:HA3	1.95	0.47
1:A:70:LEU:O	1:A:74:ILE:HG13	2.14	0.47
3:C:86:VAL:O	3:C:87:LEU:C	2.57	0.47
3:C:104:VAL:HG22	3:C:147:LEU:HD21	1.95	0.47
4:D:108:THR:O	4:D:112:GLN:HG3	2.14	0.47
2:B:12:ARG:NH1	4:D:121:VAL:HG12	2.29	0.47
3:C:99:GLN:C	3:C:101:ARG:H	2.21	0.47
4:D:14:GLU:HA	4:D:17:GLN:HB2	1.97	0.47
4:D:58:ASN:OD1	4:D:60:GLU:HB2	2.15	0.47
4:D:90:SER:O	4:D:93:ASN:HB3	2.15	0.47
2:B:5:LEU:O	2:B:8:LEU:HB2	2.15	0.47
2:B:37:ALA:O	2:B:40:PRO:HD3	2.15	0.47
3:C:7:GLU:CD	3:C:7:GLU:H	2.22	0.47
4:D:15:TRP:C	4:D:17:GLN:H	2.23	0.47
2:B:127:ASP:O	2:B:131:SER:HB2	2.15	0.46
3:C:78:ILE:O	3:C:81:LEU:HB2	2.14	0.46
1:A:13:GLN:NE2	3:C:125:SER:HB3	2.30	0.46
3:C:104:VAL:HA	3:C:108:HIS:ND1	2.30	0.46
1:A:45:LEU:HD21	1:A:99:LEU:CD2	2.45	0.46
1:A:76:GLN:HE21	1:A:79:GLN:HG3	1.81	0.46
1:A:89:HIS:O	1:A:93:GLN:HG3	2.16	0.46
4:D:32:GLY:O	4:D:35:ILE:HB	2.15	0.46
2:B:27:LEU:HD12	2:B:114:ALA:HB1	1.97	0.46
3:C:43:ASP:OD1	3:C:46:ALA:HB2	2.15	0.46
4:D:17:GLN:NE2	4:D:122:LEU:HD11	2.30	0.46
3:C:84:PRO:O	3:C:85:ALA:C	2.59	0.46
2:B:43:ARG:HD3	2:B:52:ASN:ND2	2.29	0.46
4:D:5:ARG:O	4:D:9:GLU:HG3	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:42:ARG:HG3	3:C:111:VAL:HG13	1.98	0.46
1:A:100:ASP:H	1:A:103:ASN:HD22	1.59	0.46
3:C:93:HIS:O	3:C:94:LEU:C	2.59	0.46
2:B:104:TYR:HB2	2:B:105:PRO:HD3	1.98	0.46
3:C:110:ASP:O	3:C:114:GLU:HG2	2.15	0.46
2:B:31:ILE:O	2:B:31:ILE:HD12	2.16	0.45
4:D:43:SER:C	4:D:45:VAL:H	2.24	0.45
2:B:89:ARG:HH11	2:B:89:ARG:HB2	1.79	0.45
3:C:19:ALA:HB1	3:C:123:VAL:CG1	2.36	0.45
1:A:8:ILE:HG21	3:C:123:VAL:HG13	1.99	0.45
4:D:106:ALA:O	4:D:109:HIS:HB2	2.16	0.45
2:B:12:ARG:O	2:B:15:ALA:HB3	2.17	0.45
2:B:39:MET:HG2	2:B:42:ALA:HB2	1.98	0.45
1:A:129:GLU:O	1:A:132:GLU:HB3	2.16	0.45
3:C:130:ASP:HB3	3:C:134:ARG:NH2	2.32	0.45
4:D:62:LEU:O	4:D:63:ALA:C	2.60	0.44
2:B:5:LEU:HD22	4:D:121:VAL:HG22	1.99	0.44
4:D:50:PHE:N	4:D:50:PHE:CD1	2.85	0.44
1:A:31:LEU:HA	1:A:114:VAL:HG21	1.99	0.44
1:A:34:GLU:O	1:A:35:LEU:C	2.60	0.44
1:A:46:PHE:O	1:A:51:VAL:HG22	2.18	0.44
4:D:43:SER:HB2	4:D:45:VAL:HG23	2.00	0.44
4:D:45:VAL:HG12	4:D:46:SER:N	2.32	0.44
3:C:18:ALA:C	3:C:20:TRP:N	2.76	0.44
1:A:17:SER:O	1:A:21:ALA:HB2	2.17	0.44
4:D:15:TRP:C	4:D:17:GLN:N	2.76	0.44
2:B:43:ARG:HH11	2:B:43:ARG:CB	2.30	0.44
4:D:142:GLY:HA2	4:D:145:LEU:HD12	2.00	0.44
2:B:49:VAL:HG23	2:B:59:PHE:HD1	1.82	0.44
1:A:3:ASN:OD1	1:A:6:GLU:HG3	2.17	0.44
2:B:53:ASP:O	2:B:55:ASP:N	2.51	0.43
4:D:54:GLY:O	4:D:61:PHE:CD2	2.71	0.43
1:A:8:ILE:O	1:A:9:LEU:C	2.59	0.43
1:A:57:ALA:HB1	4:D:84:THR:HA	2.00	0.43
4:D:35:ILE:HA	4:D:117:VAL:HG21	2.00	0.43
1:A:100:ASP:N	1:A:103:ASN:HD22	2.16	0.43
2:B:97:ARG:H	2:B:97:ARG:CD	2.26	0.43
3:C:105:ALA:O	3:C:106:ALA:C	2.61	0.43
2:B:125:ASP:OD1	2:B:128:ALA:HB2	2.19	0.43
2:B:104:TYR:N	2:B:105:PRO:CD	2.82	0.43
1:A:99:LEU:HD22	5:A:200:HEM:HAC	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:121:PRO:HG3	3:C:127:PHE:CD2	2.54	0.43
2:B:97:ARG:HD2	2:B:97:ARG:N	2.30	0.43
2:B:108:VAL:O	2:B:112:ILE:HG13	2.19	0.43
1:A:111:VAL:O	1:A:115:VAL:HG23	2.19	0.42
2:B:31:ILE:HG13	2:B:32:TRP:N	2.33	0.42
1:A:90:LEU:O	1:A:93:GLN:N	2.49	0.42
4:D:118:VAL:HG11	4:D:130:TRP:CZ2	2.54	0.42
1:A:36:PHE:CE2	1:A:43:ARG:HA	2.54	0.42
3:C:42:ARG:HH11	3:C:114:GLU:CD	2.28	0.42
3:C:18:ALA:C	3:C:20:TRP:H	2.26	0.42
3:C:31:LEU:HD13	3:C:123:VAL:HG21	2.02	0.42
3:C:144:SER:O	3:C:147:LEU:HB3	2.20	0.42
4:D:36:PHE:CZ	4:D:68:VAL:HG21	2.51	0.42
4:D:89:LEU:HD21	4:D:137:ILE:HG23	2.02	0.42
1:A:3:ASN:O	1:A:7:GLN:HG3	2.19	0.42
1:A:73:LEU:HD23	1:A:134:ILE:HG23	2.01	0.42
1:A:79:GLN:O	1:A:83:ILE:HG22	2.19	0.42
4:D:113:ALA:O	4:D:117:VAL:HG23	2.19	0.42
2:B:36:PHE:HD1	2:B:54:ILE:HD11	1.84	0.42
4:D:76:ILE:C	4:D:78:SER:H	2.27	0.42
1:A:8:ILE:CD1	3:C:123:VAL:HG13	2.48	0.41
2:B:1:ASP:C	2:B:131:SER:OG	2.63	0.41
3:C:25:SER:O	3:C:29:VAL:HG23	2.20	0.41
3:C:32:ALA:HB2	3:C:71:VAL:HG21	2.02	0.41
1:A:134:ILE:O	1:A:138:ILE:HG13	2.19	0.41
3:C:20:TRP:CD1	3:C:20:TRP:C	2.97	0.41
3:C:43:ASP:CG	3:C:46:ALA:HB2	2.45	0.41
1:A:57:ALA:HA	4:D:84:THR:HG22	2.03	0.41
2:B:3:THR:H	2:B:6:ASN:ND2	2.18	0.41
4:D:108:THR:C	4:D:111:SER:HG	2.23	0.41
1:A:51:VAL:O	1:A:52:ASP:C	2.64	0.41
4:D:54:GLY:O	4:D:55:ASP:C	2.62	0.41
4:D:121:VAL:O	4:D:123:PRO:HD3	2.20	0.41
1:A:23:ASN:HD21	1:A:26:ALA:HB2	1.85	0.41
1:A:131:TYR:HE1	5:A:200:HEM:HBB1	1.84	0.41
3:C:91:LEU:CD2	3:C:139:ILE:HG23	2.48	0.41
3:C:99:GLN:C	3:C:101:ARG:N	2.79	0.41
2:B:35:LEU:CD1	2:B:39:MET:HE2	2.47	0.41
3:C:40:PHE:CE1	3:C:47:GLN:HA	2.55	0.41
3:C:125:SER:O	3:C:126:CYS:SG	2.78	0.41
1:A:14:TRP:NE1	1:A:71:ASN:ND2	2.46	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:79:ASP:OD2	3:C:64:ARG:NH2	2.53	0.41
1:A:100:ASP:OD1	1:A:102:SER:N	2.54	0.41
2:B:140:ILE:HD12	5:B:200:HEM:HBB2	2.03	0.41
2:B:36:PHE:HB3	2:B:43:ARG:NH1	2.36	0.40
3:C:60:SER:C	3:C:64:ARG:HD3	2.46	0.40
3:C:82:ASN:HD22	3:C:82:ASN:HA	1.65	0.40
2:B:58:GLU:H	2:B:58:GLU:CD	2.29	0.40
4:D:27:SER:C	4:D:29:SER:N	2.79	0.40
4:D:97:ARG:HA	4:D:145:LEU:HD21	2.03	0.40
1:A:34:GLU:O	1:A:37:ASN:N	2.54	0.40
2:B:132:CYS:O	2:B:136:ILE:HG13	2.21	0.40
2:B:27:LEU:CD1	2:B:114:ALA:HB1	2.52	0.40
3:C:36:PHE:O	3:C:37:ALA:C	2.63	0.40
3:C:125:SER:C	3:C:126:CYS:SG	3.05	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all 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	138/140 (99%)	122 (88%)	10 (7%)	6 (4%)	2	4
2	B	140/142 (99%)	109 (78%)	27 (19%)	4 (3%)	3	8
3	C	144/147 (98%)	120 (83%)	18 (12%)	6 (4%)	2	3
4	D	143/145 (99%)	118 (82%)	21 (15%)	4 (3%)	4	9
All	All	565/574 (98%)	469 (83%)	76 (14%)	20 (4%)	3	6

All (20) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	49	VAL

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Mol	Chain	Res	Type
3	C	104	VAL
2	B	18	TYR
2	B	54	ILE
3	C	54	SER
4	D	22	ALA
1	A	21	ALA
1	A	22	GLU
2	B	19	GLY
3	C	43	ASP
1	A	124	GLN
1	A	125	HIS
3	C	59	ASP
4	D	2	CYS
1	A	139	SER
2	B	102	ALA
3	C	81	LEU
3	C	93	HIS
4	D	125	ALA
4	D	127	ILE

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	110/110 (100%)	109 (99%)	1 (1%)	70	84
2	B	113/113 (100%)	108 (96%)	5 (4%)	25	49
3	C	110/111 (99%)	105 (96%)	5 (4%)	24	48
4	D	108/108 (100%)	105 (97%)	3 (3%)	38	62
All	All	441/442 (100%)	427 (97%)	14 (3%)	34	59

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

Mol	Chain	Res	Type
1	A	34	GLU
2	B	26	LEU

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Mol	Chain	Res	Type
2	B	31	ILE
2	B	78	ASP
2	B	89	ARG
2	B	95	ASP
3	C	7	GLU
3	C	74	LEU
3	C	81	LEU
3	C	89	GLU
3	C	134	ARG
4	D	28	GLU
4	D	110	LEU
4	D	128	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (29) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	16	GLN
1	A	37	ASN
1	A	39	GLN
1	A	62	HIS
1	A	71	ASN
1	A	76	GLN
1	A	103	ASN
2	B	6	ASN
2	B	13	GLN
2	B	23	ASN
2	B	50	ASN
2	B	52	ASN
2	B	84	ASN
2	B	93	GLN
2	B	134	ASN
3	C	11	ASN
3	C	33	GLN
3	C	47	GLN
3	C	57	ASN
3	C	66	HIS
3	C	79	ASN
3	C	82	ASN
3	C	90	GLN
3	C	99	GLN
3	C	122	GLN
3	C	133	ASN

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Mol	Chain	Res	Type
3	C	141	ASN
4	D	17	GLN
4	D	93	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

8 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
6	OXY	C	2201	5	1,1,1	0.00	0	-		
6	OXY	A	201	5	1,1,1	0.01	0	-		
6	OXY	D	3201	5	1,1,1	0.00	0	-		
5	HEM	A	200	1,6	50,50,50	1.32	5 (10%)	67,82,82	0.74	2 (2%)
5	HEM	D	200	6,4	50,50,50	1.37	4 (8%)	67,82,82	0.75	2 (2%)
6	OXY	B	1201	5	1,1,1	0.00	0	-		
5	HEM	B	200	6,2	50,50,50	1.31	5 (10%)	67,82,82	0.73	1 (1%)
5	HEM	C	200	3,6	50,50,50	1.34	4 (8%)	67,82,82	0.72	1 (1%)

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
5	HEM	D	200	6,4	-	9/14/54/54	-
5	HEM	A	200	1,6	-	9/14/54/54	-
5	HEM	B	200	6,2	-	8/14/54/54	-
5	HEM	C	200	3,6	-	9/14/54/54	-

All (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	200	HEM	FE-NB	4.31	2.08	1.94
5	A	200	HEM	FE-NC	4.26	2.09	1.95
5	C	200	HEM	FE-NC	4.19	2.09	1.95
5	D	200	HEM	FE-NB	4.19	2.07	1.94
5	B	200	HEM	FE-NC	4.18	2.09	1.95
5	B	200	HEM	FE-NB	4.12	2.07	1.94
5	D	200	HEM	FE-NC	4.07	2.08	1.95
5	C	200	HEM	FE-NB	3.82	2.06	1.94
5	D	200	HEM	FE-ND	3.55	2.05	1.94
5	D	200	HEM	FE-NA	3.52	2.06	1.95
5	C	200	HEM	FE-ND	3.36	2.05	1.94
5	C	200	HEM	FE-NA	3.22	2.05	1.95
5	B	200	HEM	FE-ND	3.18	2.04	1.94
5	A	200	HEM	FE-NA	2.74	2.04	1.95
5	B	200	HEM	FE-NA	2.73	2.04	1.95
5	A	200	HEM	FE-ND	2.62	2.03	1.94
5	A	200	HEM	C4D-ND	-2.18	1.36	1.40
5	B	200	HEM	C4D-ND	-2.07	1.36	1.40

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	200	HEM	C4D-ND-C1D	2.38	108.03	105.21
5	D	200	HEM	C3B-C2B-C1B	-2.24	104.73	106.41
5	C	200	HEM	CAD-CBD-CGD	-2.12	108.04	113.67
5	B	200	HEM	CAD-CBD-CGD	-2.07	108.17	113.67
5	A	200	HEM	CAD-CBD-CGD	-2.05	108.23	113.67
5	D	200	HEM	C4D-ND-C1D	2.03	107.61	105.21

There are no chirality outliers.

All (35) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	200	HEM	C2B-C3B-CAB-CBB
5	A	200	HEM	C4B-C3B-CAB-CBB
5	B	200	HEM	C2B-C3B-CAB-CBB
5	B	200	HEM	C2C-C3C-CAC-CBC
5	C	200	HEM	C2B-C3B-CAB-CBB
5	C	200	HEM	C4B-C3B-CAB-CBB
5	D	200	HEM	C2B-C3B-CAB-CBB
5	D	200	HEM	C2C-C3C-CAC-CBC
5	D	200	HEM	C2A-CAA-CBA-CGA
5	C	200	HEM	C2A-CAA-CBA-CGA
5	C	200	HEM	C2C-C3C-CAC-CBC
5	B	200	HEM	C4B-C3B-CAB-CBB
5	B	200	HEM	C4C-C3C-CAC-CBC
5	D	200	HEM	C4B-C3B-CAB-CBB
5	D	200	HEM	C4C-C3C-CAC-CBC
5	A	200	HEM	C3D-CAD-CBD-CGD
5	A	200	HEM	C2C-C3C-CAC-CBC
5	C	200	HEM	C4C-C3C-CAC-CBC
5	C	200	HEM	CAD-CBD-CGD-O1D
5	C	200	HEM	CAA-CBA-CGA-O2A
5	A	200	HEM	CAA-CBA-CGA-O2A
5	A	200	HEM	CAD-CBD-CGD-O1D
5	B	200	HEM	CAD-CBD-CGD-O1D
5	C	200	HEM	CAA-CBA-CGA-O1A
5	A	200	HEM	CAD-CBD-CGD-O2D
5	B	200	HEM	CAD-CBD-CGD-O2D
5	A	200	HEM	CAA-CBA-CGA-O1A
5	C	200	HEM	CAD-CBD-CGD-O2D
5	D	200	HEM	CAA-CBA-CGA-O1A
5	D	200	HEM	CAA-CBA-CGA-O2A
5	B	200	HEM	CAA-CBA-CGA-O2A
5	D	200	HEM	CAD-CBD-CGD-O1D
5	D	200	HEM	CAD-CBD-CGD-O2D
5	B	200	HEM	CAA-CBA-CGA-O1A
5	A	200	HEM	C4C-C3C-CAC-CBC

There are no ring outliers.

3 monomers are involved in 10 short contacts:

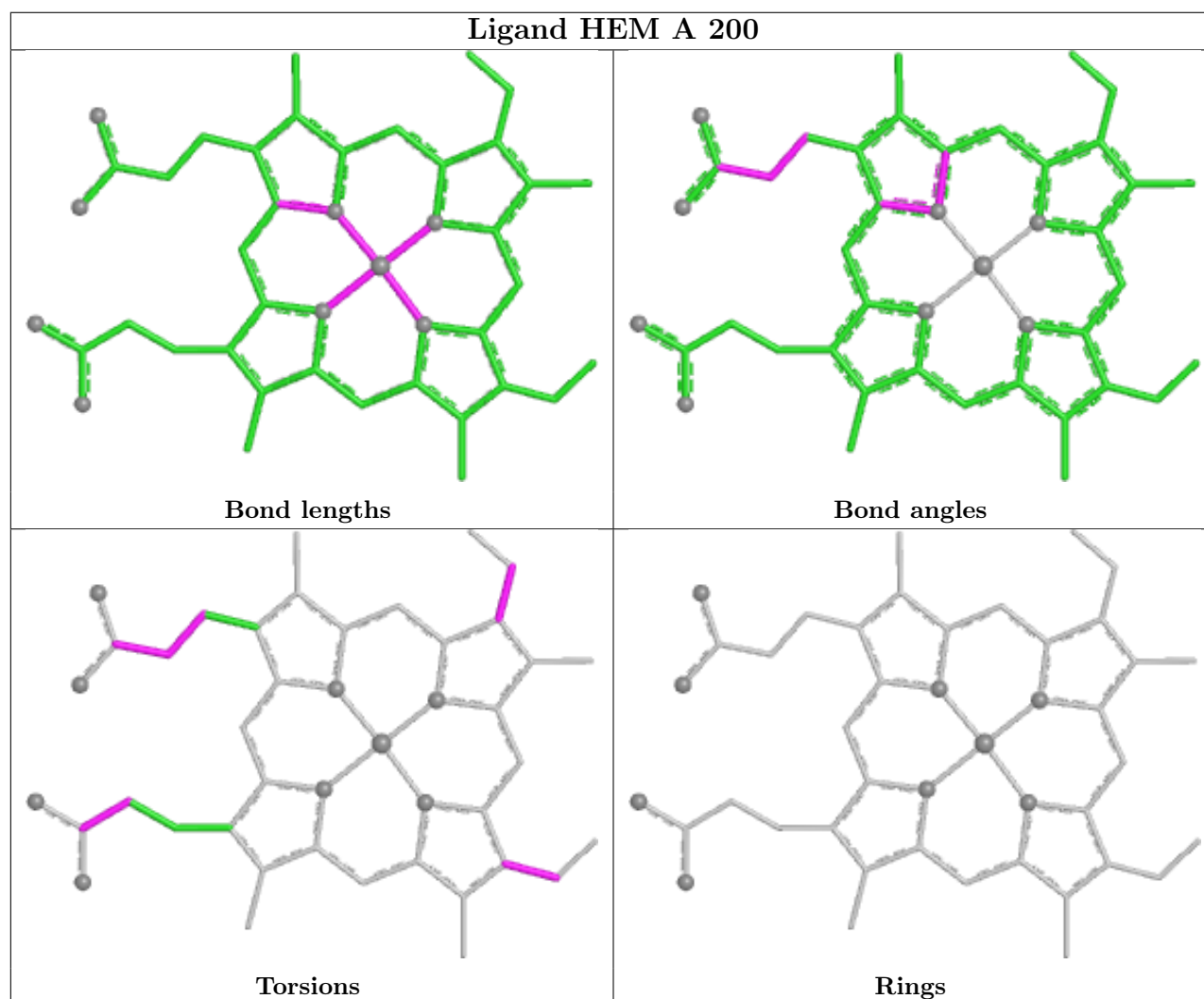
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	200	HEM	4	0

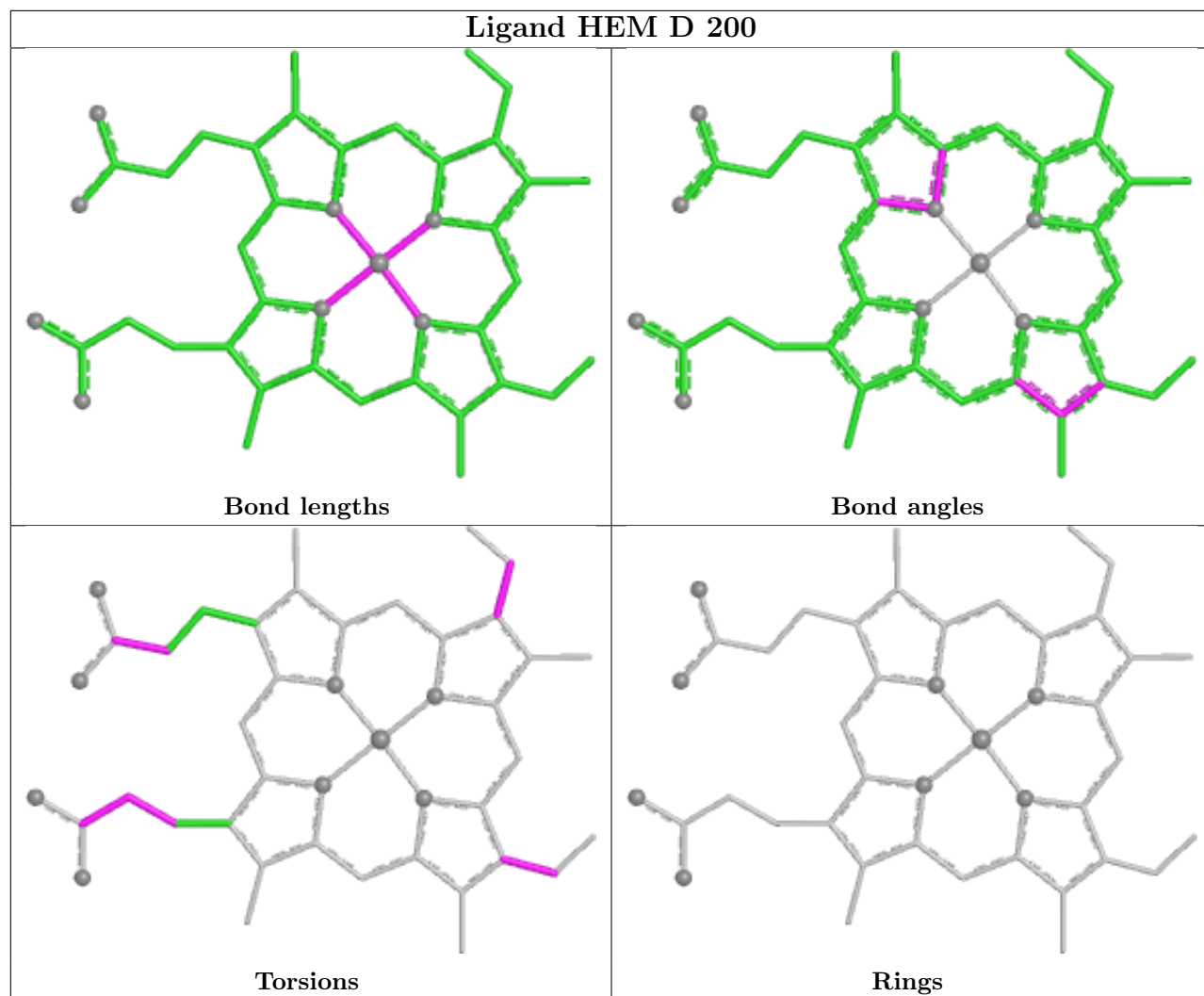
Continued on next page...

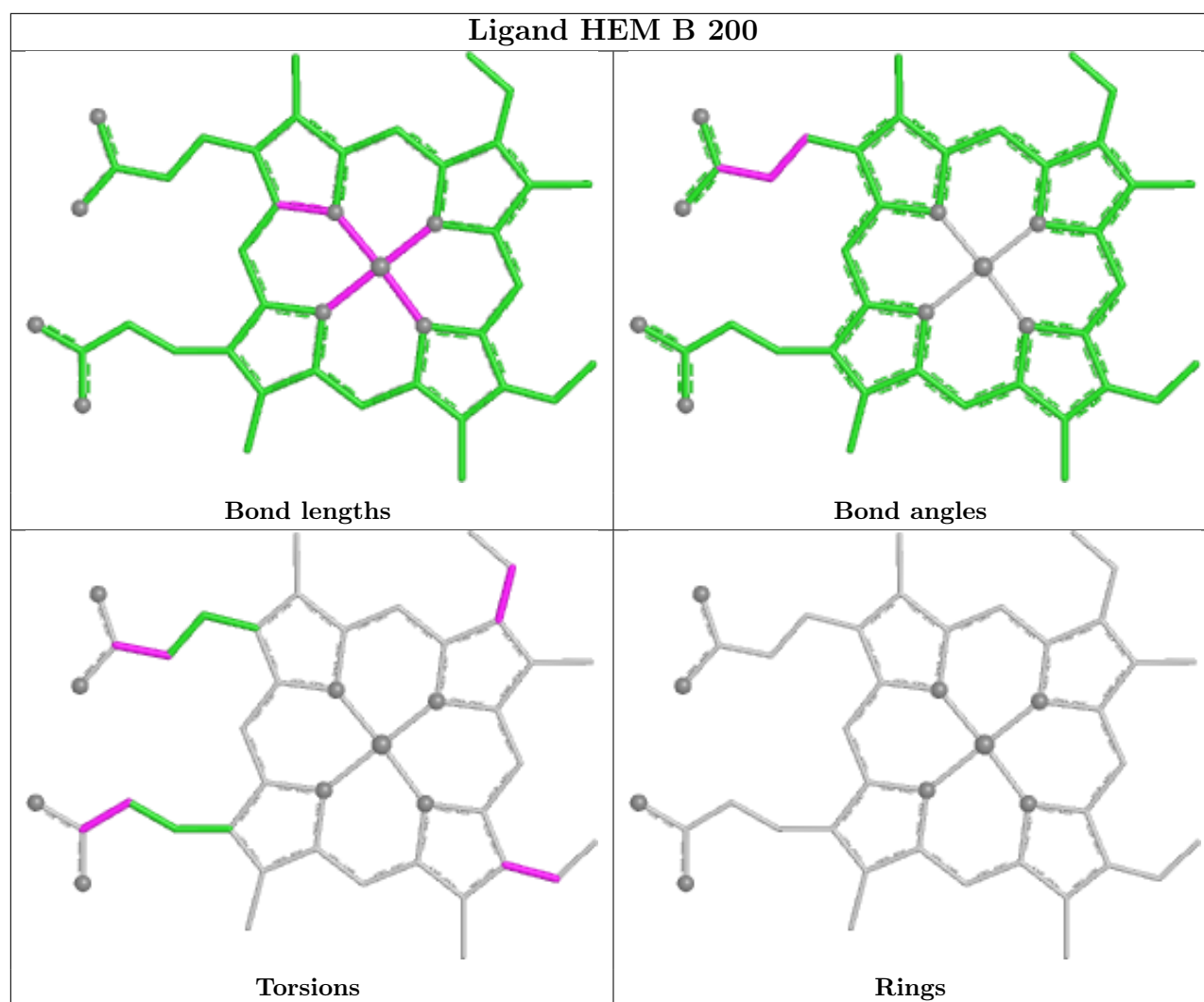
Continued from previous page...

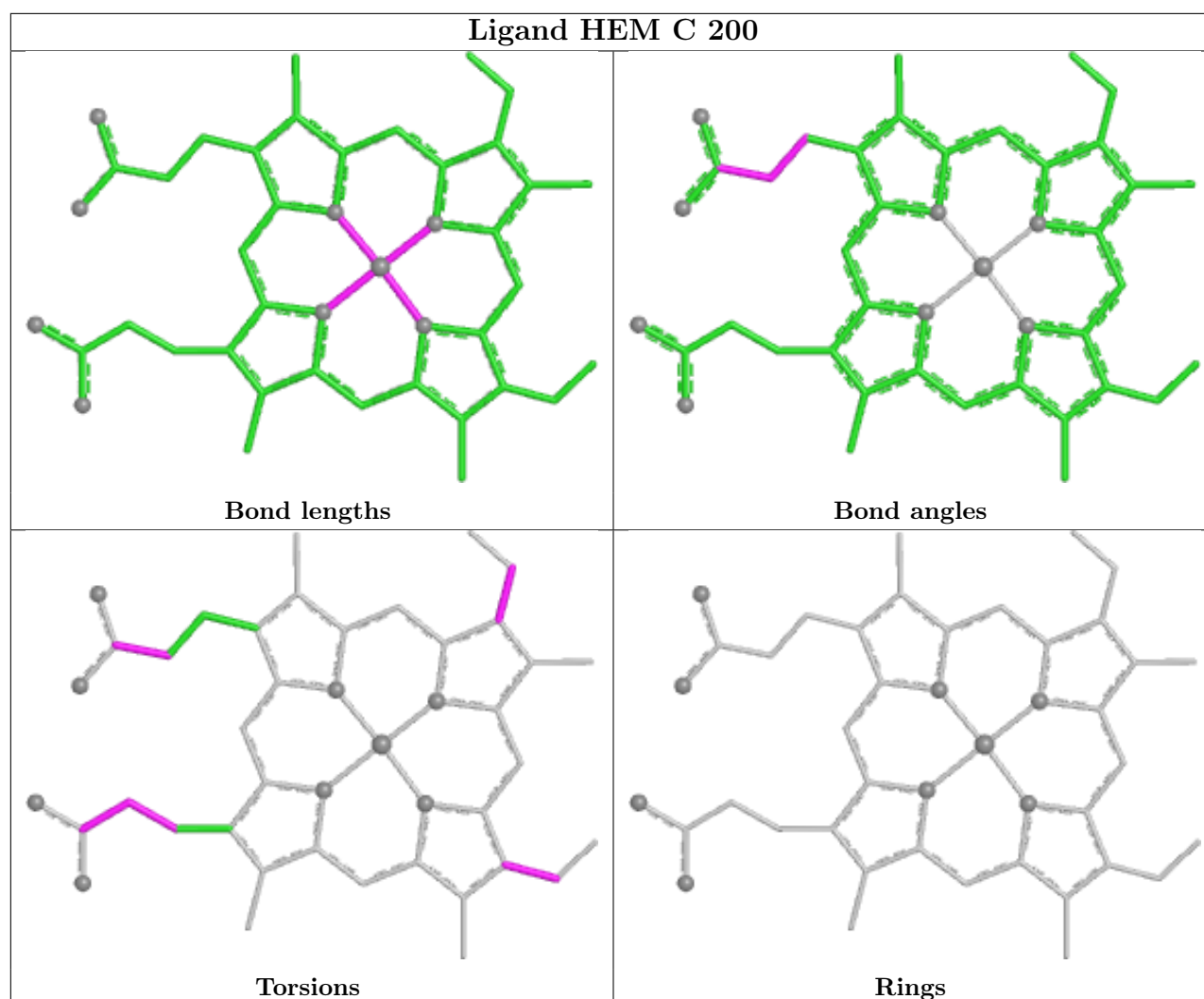
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	D	200	HEM	2	0
5	B	200	HEM	4	0

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.

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.