



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 28, 2026 – 09:00 PM UTC

PDB ID : 4NOS / pdb_00004nos
Title : HUMAN INDUCIBLE NITRIC OXIDE SYNTHASE WITH INHIBITOR
Authors : Fischmann, T.O.; Weber, P.C.
Deposited on : 1999-02-03
Resolution : 2.25 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)	:	2.49

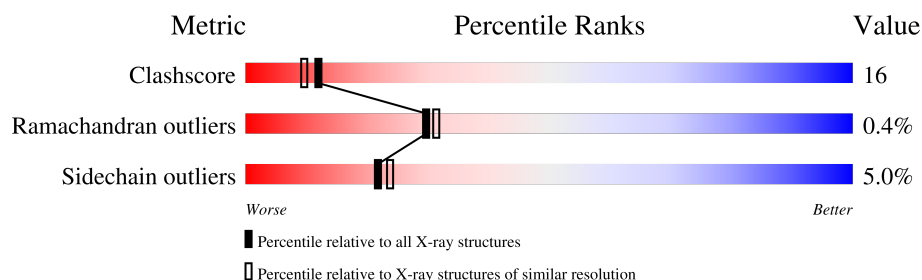
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.25 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	2005 (2.26-2.26)
Ramachandran outliers	187476	1965 (2.26-2.26)
Sidechain outliers	187428	1966 (2.26-2.26)

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	427	 66% 29% . .
1	B	427	 60% 33% 5% .
1	C	427	 64% 30% . .
1	D	427	 69% 26% . .

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
5	ITU	B	1012	-	-	X	-

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 15317 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 INDUCIBLE NITRIC OXIDE SYNTHASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	421	Total	C	N	O	S	0	0	1
			3420	2186	600	612	22			
1	B	421	Total	C	N	O	S	0	0	1
			3420	2186	600	612	22			
1	C	421	Total	C	N	O	S	0	0	1
			3420	2186	600	612	22			
1	D	421	Total	C	N	O	S	0	0	1
			3420	2186	600	612	22			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	423	ILE	LEU	conflict	UNP P35228
B	423	ILE	LEU	conflict	UNP P35228
C	423	ILE	LEU	conflict	UNP P35228
D	423	ILE	LEU	conflict	UNP P35228

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

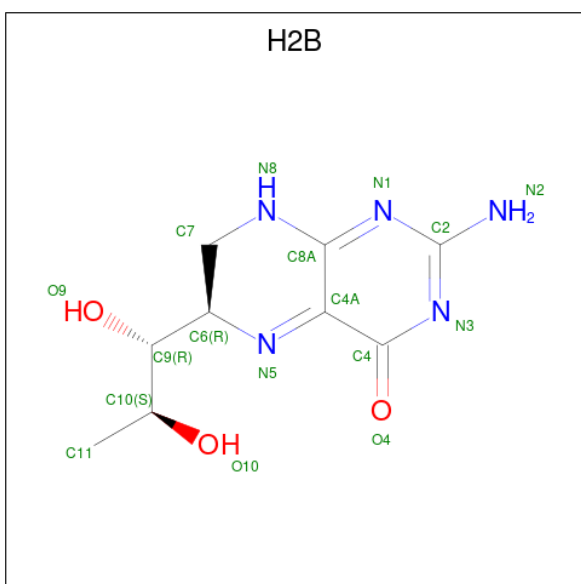
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Zn	0	0
			1	1		
2	C	1	Total	Zn	0	0
			1	1		

- Molecule 3 is PROTOPORPHYRIN IX CONTAINING FE (CCD ID: HEM) (formula: C₃₄H₃₂FeN₄O₄).



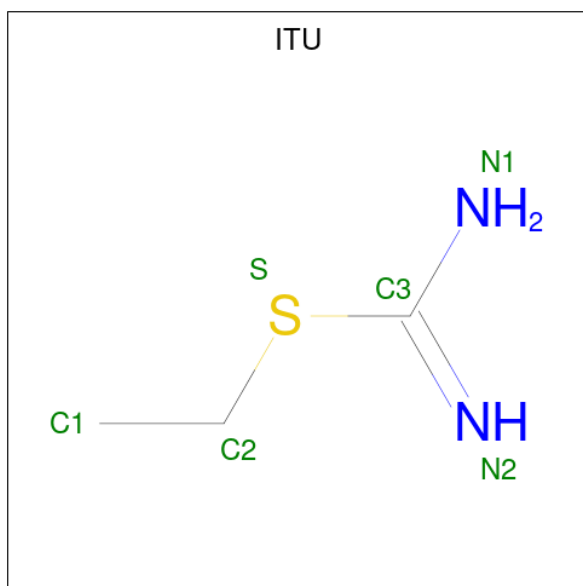
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
3	B	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
3	C	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
3	D	1	Total 43	C 34	Fe 1	N 4	O 4	0	0

- Molecule 4 is 2-AMINO-6-(1,2-DIHYDROXY-PROPYL)-7,8-DIHYDRO-6H-PTERIDIN-4-ONE (CCD ID: H2B) (formula: $C_9H_{13}N_5O_3$).



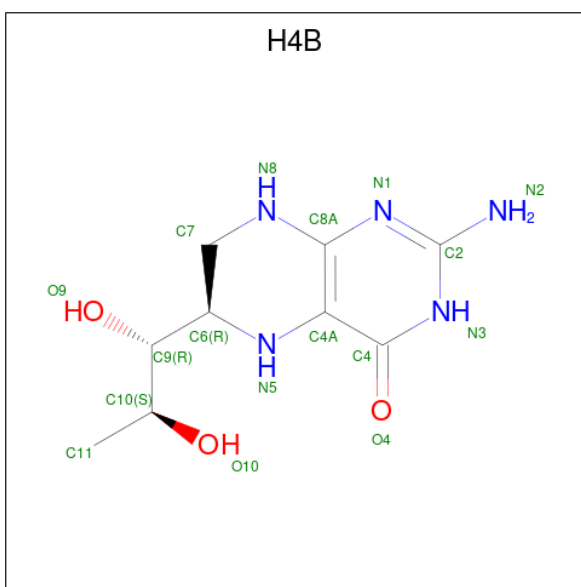
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			17	9	5	3		

- Molecule 5 is ETHYLISOTHIOUREA (CCD ID: ITU) (formula: $C_3H_8N_2S$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	S	0	0
			6	3	2	1		
5	B	1	Total	C	N	S	0	0
			6	3	2	1		
5	C	1	Total	C	N	S	0	0
			6	3	2	1		
5	D	1	Total	C	N	S	0	0
			6	3	2	1		

- Molecule 6 is 5,6,7,8-TETRAHYDROBIOPTERIN (CCD ID: H4B) (formula: $C_9H_{15}N_5O_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	B	1	Total	C	N	O	0	0
			17	9	5	3		
6	C	1	Total	C	N	O	0	0
			17	9	5	3		
6	D	1	Total	C	N	O	0	0
			17	9	5	3		

- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	336	Total	O	0	0
			336	336		
7	B	335	Total	O	0	0
			335	335		
7	C	339	Total	O	0	0
			339	339		
7	D	361	Total	O	0	0
			361	361		

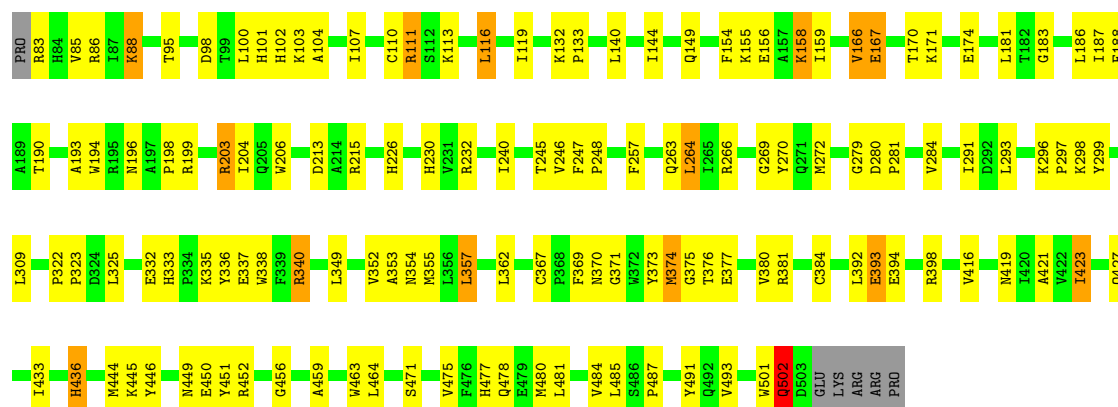
3 Residue-property plots

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.

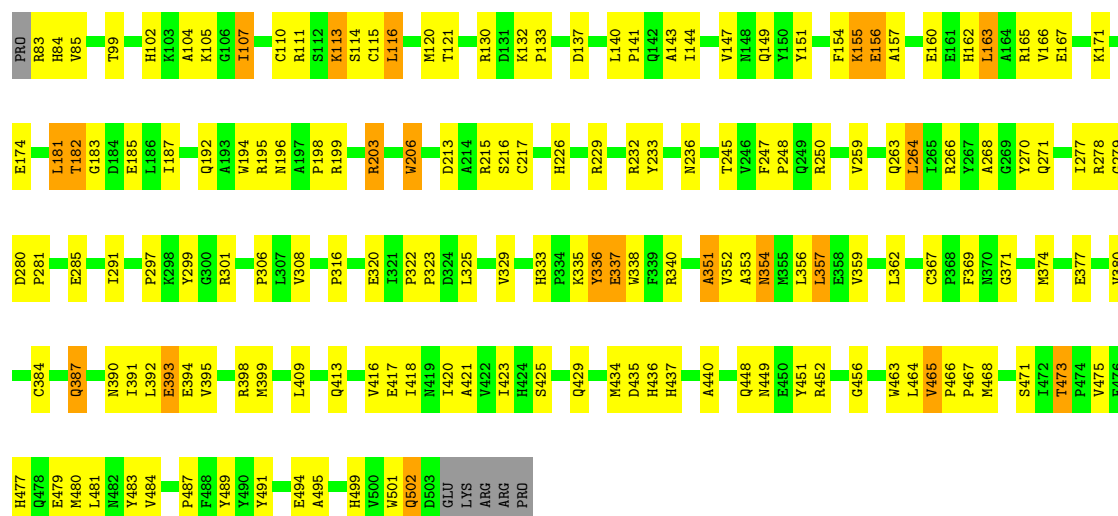
• Molecule 1: INDUCIBLE NITRIC OXIDE SYNTHASE

Chain A: 



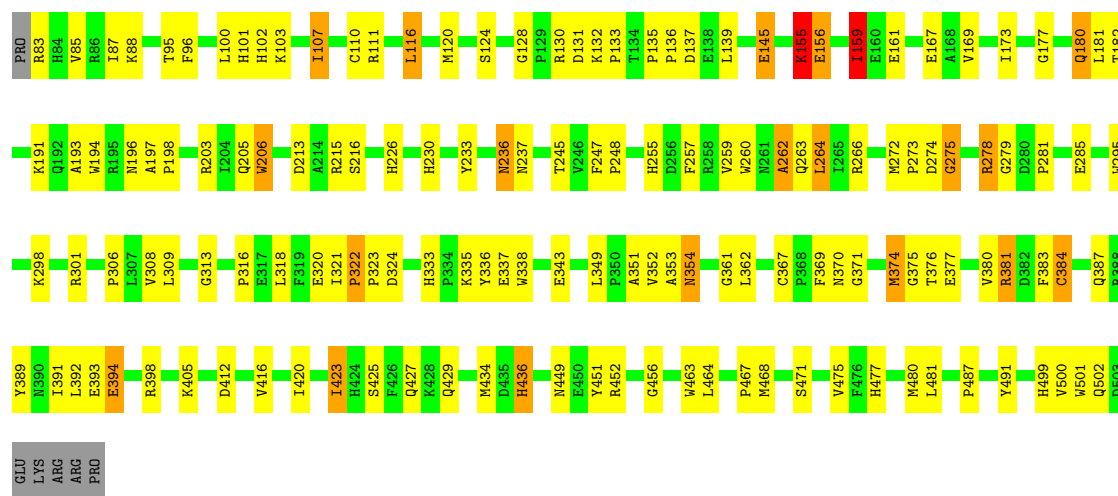
• Molecule 1: INDUCIBLE NITRIC OXIDE SYNTHASE

Chain B: 



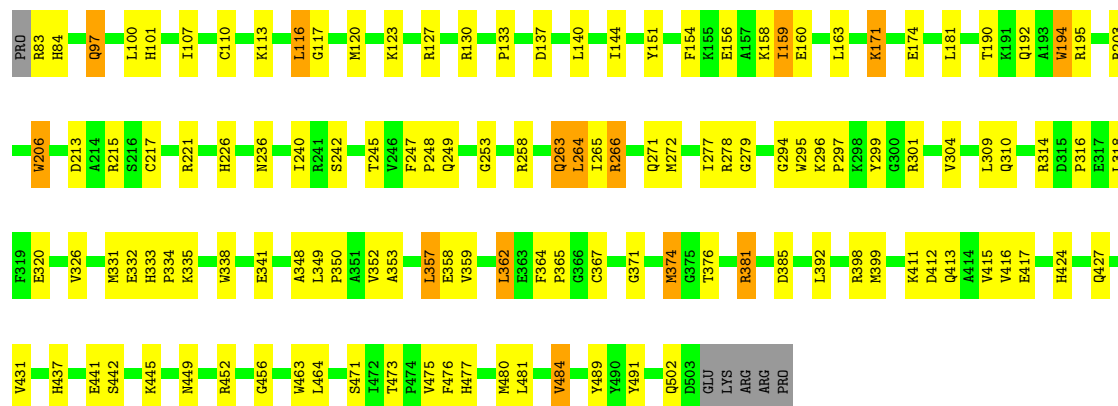
• Molecule 1: INDUCIBLE NITRIC OXIDE SYNTHASE

Chain C:  64% 30% . .



• Molecule 1: INDUCIBLE NITRIC OXIDE SYNTHASE

Chain D:  69% 26% . .



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	89.76Å 156.67Å 190.04Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	8.00 – 2.25	Depositor
% Data completeness (in resolution range)	88.7 (8.00-2.25)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.199 , 0.289	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	15317	wwPDB-VP
Average B, all atoms (Å ²)	26.0	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: ZN, ITU, H4B, HEM, H2B

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	0.63	1/3519 (0.0%)	1.03	16/4776 (0.3%)
1	B	0.63	1/3519 (0.0%)	1.06	27/4776 (0.6%)
1	C	0.64	1/3519 (0.0%)	1.08	21/4776 (0.4%)
1	D	0.65	1/3519 (0.0%)	1.06	16/4776 (0.3%)
All	All	0.64	4/14076 (0.0%)	1.06	80/19104 (0.4%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	502	GLN	C-N	-5.76	1.25	1.33
1	C	502	GLN	C-N	-5.70	1.25	1.33
1	B	502	GLN	C-N	-5.67	1.25	1.33
1	A	502	GLN	C-N	-5.57	1.25	1.33

All (80) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	128	GLY	CA-C-N	8.89	128.94	119.78
1	C	128	GLY	C-N-CA	8.89	128.94	119.78
1	C	374	MET	N-CA-C	-8.25	95.78	109.07
1	B	374	MET	N-CA-C	-7.31	96.60	108.52
1	B	499	HIS	N-CA-C	7.16	120.63	109.96
1	B	203	ARG	N-CA-C	7.11	121.40	112.87
1	D	381	ARG	N-CA-C	6.92	120.16	111.24
1	C	375	GLY	N-CA-C	6.86	121.20	112.83
1	C	155	LYS	N-CA-C	-6.86	104.17	112.54
1	C	354	ASN	N-CA-C	6.82	120.50	112.72
1	B	465	VAL	CA-C-N	6.78	124.61	119.66
1	B	465	VAL	C-N-CA	6.78	124.61	119.66
1	D	120	MET	N-CA-C	6.70	118.46	111.03

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	374	MET	N-CA-C	-6.68	97.84	108.73
1	D	374	MET	N-CA-C	-6.63	98.43	109.24
1	B	120	MET	N-CA-C	6.59	118.16	110.97
1	D	194	TRP	N-CA-C	-6.52	104.10	111.14
1	A	98	ASP	N-CA-C	6.47	119.11	109.59
1	A	493	VAL	N-CA-C	-6.41	101.29	109.30
1	C	120	MET	N-CA-C	6.32	117.83	111.07
1	A	354	ASN	N-CA-C	6.31	120.76	111.87
1	D	371	GLY	N-CA-C	-6.27	98.32	113.18
1	C	259	VAL	N-CA-C	-6.25	98.86	107.99
1	D	392	LEU	N-CA-C	6.22	118.61	111.02
1	B	107	ILE	N-CA-C	6.20	119.60	113.53
1	C	182	THR	N-CA-C	-6.20	102.35	110.53
1	A	373	TYR	N-CA-C	6.18	119.96	110.20
1	C	371	GLY	N-CA-C	-6.03	98.89	113.18
1	D	331	MET	N-CA-C	5.99	117.96	108.67
1	D	491	TYR	N-CA-C	-5.93	101.16	110.36
1	A	436	HIS	N-CA-C	5.92	118.69	111.82
1	B	491	TYR	N-CA-C	-5.89	102.94	110.53
1	A	155	LYS	N-CA-C	-5.84	105.00	111.36
1	C	384	CYS	N-CA-C	5.84	121.44	113.97
1	C	203	ARG	N-CA-C	5.80	119.83	112.87
1	D	484	VAL	N-CA-C	5.79	115.94	107.37
1	A	203	ARG	N-CA-C	5.78	123.11	110.80
1	D	159	ILE	N-CA-C	5.72	115.91	110.42
1	A	166	VAL	N-CA-C	-5.71	105.16	110.53
1	C	436	HIS	N-CA-C	5.70	118.26	111.71
1	D	240	ILE	N-CA-C	5.64	117.57	109.51
1	A	309	LEU	N-CA-C	5.63	118.13	109.07
1	B	354	ASN	N-CA-C	5.61	119.44	112.59
1	C	491	TYR	N-CA-C	-5.61	102.61	110.35
1	C	262	ALA	N-CA-C	-5.58	105.28	111.36
1	D	473	THR	CA-C-N	5.58	125.41	119.28
1	D	473	THR	C-N-CA	5.58	125.41	119.28
1	B	467	PRO	N-CA-C	-5.57	108.29	114.92
1	A	371	GLY	N-CA-C	-5.52	99.37	110.69
1	B	157	ALA	N-CA-C	5.52	118.32	110.10
1	B	495	ALA	N-CA-C	5.50	117.72	111.11
1	D	203	ARG	N-CA-C	5.48	119.45	112.87
1	C	322	PRO	CA-C-N	5.46	125.80	119.47
1	C	322	PRO	C-N-CA	5.46	125.80	119.47
1	C	420	ILE	N-CA-C	-5.39	105.27	110.72

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	159	ILE	N-CA-C	5.37	116.74	110.62
1	A	280	ASP	CA-C-N	5.37	125.19	119.28
1	A	280	ASP	C-N-CA	5.37	125.19	119.28
1	B	114	SER	N-CA-C	5.34	116.53	108.46
1	B	182	THR	N-CA-C	-5.34	103.85	110.41
1	B	479	GLU	N-CA-C	-5.34	102.98	110.50
1	A	240	ILE	N-CA-C	5.30	117.09	109.51
1	B	259	VAL	N-CA-C	-5.29	100.27	107.99
1	C	145	GLU	N-CA-C	-5.29	105.43	111.14
1	B	121	THR	CA-C-N	-5.28	114.52	119.85
1	B	121	THR	C-N-CA	-5.28	114.52	119.85
1	C	107	ILE	N-CA-C	5.23	118.17	113.10
1	B	473	THR	CA-C-N	5.22	125.02	119.28
1	B	473	THR	C-N-CA	5.22	125.02	119.28
1	B	351	ALA	N-CA-C	5.21	116.82	107.80
1	B	390	ASN	N-CA-C	5.20	118.42	111.24
1	A	491	TYR	N-CA-C	-5.20	102.77	110.46
1	B	336	TYR	N-CA-C	5.16	117.08	107.99
1	B	280	ASP	CA-C-N	5.15	124.94	119.28
1	B	280	ASP	C-N-CA	5.15	124.94	119.28
1	B	206	TRP	N-CA-C	5.11	118.28	111.75
1	D	117	GLY	N-CA-C	5.08	120.04	114.40
1	A	421	ALA	N-CA-C	-5.05	105.67	111.07
1	D	294	GLY	N-CA-C	5.04	122.06	115.36
1	B	371	GLY	N-CA-C	-5.01	101.31	113.18

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	3420	0	3326	108	0
1	B	3420	0	3326	127	0
1	C	3420	0	3326	106	1
1	D	3420	0	3326	94	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	1	0	0	0	0
2	C	1	0	0	0	0
3	A	43	0	30	9	0
3	B	43	0	30	4	0
3	C	43	0	30	8	0
3	D	43	0	30	9	0
4	A	17	0	13	0	0
5	A	6	0	7	2	0
5	B	6	0	7	4	0
5	C	6	0	7	3	0
5	D	6	0	7	2	0
6	B	17	0	15	1	0
6	C	17	0	15	2	0
6	D	17	0	15	1	0
7	A	336	0	0	10	0
7	B	335	0	0	15	1
7	C	339	0	0	16	0
7	D	361	0	0	9	0
All	All	15317	0	13510	432	1

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 16.

All (432) 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:D:217:CYS:SG	7:D:3344:HOH:O	2.28	0.90
1:C:481:LEU:HD22	1:D:116:LEU:HD22	1.52	0.90
3:C:2010:HEM:HMC2	3:C:2010:HEM:HBC2	1.56	0.88
1:A:272:MET:HE1	1:A:299:TYR:CD1	2.10	0.86
1:A:88:LYS:HB2	1:A:95:THR:HG22	1.57	0.84
3:D:3010:HEM:HBC2	3:D:3010:HEM:HMC2	1.61	0.82
3:B:1010:HEM:HMC2	3:B:1010:HEM:HBC2	1.63	0.81
1:B:155:LYS:H	1:B:155:LYS:HD3	1.45	0.80
1:A:111:ARG:NH1	1:A:113:LYS:HB3	2.00	0.77
1:A:110:CYS:HB3	1:B:110:CYS:HB3	1.66	0.77
1:B:194:TRP:CE3	1:B:206:TRP:HA	2.21	0.76
1:C:194:TRP:CE3	1:C:206:TRP:HA	2.21	0.75
1:D:213:ASP:OD1	1:D:215:ARG:HD3	1.86	0.75
1:B:335:LYS:HE3	1:B:336:TYR:CZ	2.22	0.75
1:C:110:CYS:HB3	1:D:110:CYS:HB3	1.67	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:272:MET:HE1	1:A:299:TYR:HD1	1.51	0.75
1:D:194:TRP:CE3	1:D:206:TRP:HA	2.21	0.74
1:D:314:ARG:HD2	7:D:3271:HOH:O	1.88	0.73
1:A:111:ARG:HH11	1:A:111:ARG:HB2	1.53	0.71
1:D:83:ARG:HE	1:D:84:HIS:HB3	1.55	0.71
1:D:226:HIS:HD2	7:D:3070:HOH:O	1.73	0.71
1:D:333:HIS:HD2	1:D:335:LYS:H	1.38	0.71
1:B:155:LYS:HD3	1:B:155:LYS:N	2.04	0.71
1:C:159:ILE:HD13	1:C:159:ILE:H	1.56	0.71
1:D:192:GLN:HE22	1:D:195:ARG:HH11	1.39	0.71
1:B:144:ILE:HD12	1:B:166:VAL:HG13	1.72	0.70
1:A:449:ASN:HA	1:A:452:ARG:NH1	2.07	0.70
1:C:369:PHE:HE1	5:C:2012:ITU:H11	1.57	0.69
1:D:113:LYS:HZ3	1:D:113:LYS:HB3	1.57	0.69
1:A:270:TYR:HB2	1:A:272:MET:HE3	1.74	0.69
1:A:352:VAL:HG11	1:A:355:MET:SD	2.32	0.69
1:B:268:ALA:HB2	7:B:1019:HOH:O	1.91	0.68
1:C:272:MET:HE2	1:C:278:ARG:HB2	1.74	0.68
1:C:338:TRP:CE3	1:C:398:ARG:HD3	2.29	0.67
1:B:434:MET:SD	7:B:1217:HOH:O	2.52	0.67
1:C:194:TRP:CH2	3:C:2010:HEM:HMC1	2.29	0.67
1:A:298:LYS:HE2	7:A:3088:HOH:O	1.94	0.67
1:A:471:SER:O	1:A:477:HIS:HE1	1.77	0.67
1:A:226:HIS:HD2	7:A:3057:HOH:O	1.79	0.66
1:C:213:ASP:OD1	1:C:215:ARG:HD3	1.96	0.66
1:A:481:LEU:HD22	1:B:116:LEU:HD22	1.78	0.66
1:B:333:HIS:HD2	1:B:335:LYS:H	1.45	0.65
1:D:442:SER:HA	1:D:445:LYS:HE3	1.79	0.65
1:B:156:GLU:CD	1:B:156:GLU:H	2.04	0.65
1:D:171:LYS:HA	1:D:171:LYS:HE2	1.78	0.64
1:B:213:ASP:OD1	1:B:215:ARG:HD3	1.98	0.64
1:B:264:LEU:HD13	1:B:353:ALA:HB2	1.80	0.64
1:B:329:VAL:HG13	1:B:429:GLN:HG3	1.80	0.63
1:D:140:LEU:O	1:D:144:ILE:HG12	1.98	0.63
1:D:271:GLN:HA	1:D:277:ILE:HD13	1.80	0.63
1:A:194:TRP:CE3	1:A:206:TRP:HA	2.34	0.62
1:A:213:ASP:OD1	1:A:215:ARG:HD3	2.00	0.62
1:C:226:HIS:HD2	7:C:2069:HOH:O	1.82	0.62
1:C:264:LEU:HD13	1:C:353:ALA:HB2	1.81	0.62
1:C:369:PHE:CE1	5:C:2012:ITU:H11	2.34	0.62
1:A:156:GLU:CD	1:A:156:GLU:H	2.07	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:333:HIS:HD2	1:B:336:TYR:H	1.46	0.62
1:B:105:LYS:HB2	1:B:107:ILE:HG12	1.82	0.62
1:B:183:GLY:O	1:B:187:ILE:HG13	2.00	0.62
1:D:113:LYS:HB3	1:D:113:LYS:NZ	2.14	0.62
1:D:352:VAL:HG23	5:D:3012:ITU:H11	1.81	0.61
1:A:83:ARG:O	1:A:102:HIS:HE1	1.83	0.61
1:A:263:GLN:CD	1:A:266:ARG:HD3	2.25	0.61
1:B:393:GLU:HB2	7:B:1207:HOH:O	2.01	0.61
1:B:83:ARG:O	1:B:102:HIS:HE1	1.83	0.61
1:B:392:LEU:HB2	7:B:1103:HOH:O	1.99	0.61
1:C:333:HIS:HD2	1:C:335:LYS:H	1.48	0.61
1:D:194:TRP:CZ3	1:D:206:TRP:HA	2.35	0.61
1:A:270:TYR:CB	1:A:272:MET:HE3	2.31	0.61
1:C:264:LEU:HB2	1:C:351:ALA:HB3	1.81	0.61
1:A:369:PHE:CE1	5:A:512:ITU:H11	2.35	0.61
1:B:84:HIS:HB3	1:B:99:THR:HG23	1.83	0.61
1:B:449:ASN:ND2	1:B:452:ARG:HH12	1.99	0.61
1:B:130:ARG:NH1	1:B:133:PRO:HA	2.16	0.60
1:A:101:HIS:CD2	1:B:111:ARG:HH21	2.18	0.60
1:B:369:PHE:CD1	5:B:1012:ITU:H11	2.36	0.60
1:A:419:ASN:O	1:A:423:ILE:HG12	2.00	0.60
1:B:333:HIS:CD2	1:B:335:LYS:H	2.19	0.60
1:B:369:PHE:CE1	5:B:1012:ITU:H11	2.37	0.60
1:C:416:VAL:HG22	1:C:468:MET:HE1	1.84	0.60
1:C:139:LEU:HD12	1:C:361:GLY:HA3	1.83	0.60
3:A:510:HEM:HMC2	3:A:510:HEM:HBC2	1.83	0.59
1:A:144:ILE:HD12	1:A:166:VAL:HG13	1.85	0.59
1:A:245:THR:O	1:A:367:CYS:HA	2.02	0.59
1:B:163:LEU:O	1:B:167:GLU:HG3	2.03	0.59
1:A:333:HIS:HD2	1:A:336:TYR:H	1.50	0.59
1:C:381:ARG:HH21	6:C:2011:H4B:C4	2.15	0.59
1:A:264:LEU:HD13	1:A:353:ALA:HB2	1.85	0.58
1:B:416:VAL:HG22	1:B:468:MET:HE3	1.85	0.58
1:B:151:TYR:CD1	1:B:154:PHE:HE2	2.21	0.58
1:D:309:LEU:O	1:D:316:PRO:HA	2.04	0.58
1:C:83:ARG:HH21	1:C:102:HIS:CE1	2.22	0.57
1:A:199:ARG:HD3	1:A:463:TRP:CD2	2.38	0.57
1:A:449:ASN:HA	1:A:452:ARG:HH12	1.69	0.57
1:B:195:ARG:HD2	1:B:206:TRP:CE3	2.39	0.57
1:A:291:ILE:HD11	1:A:297:PRO:HB3	1.86	0.57
1:C:263:GLN:OE1	1:C:266:ARG:HD3	2.05	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:359:VAL:O	1:D:362:LEU:HB2	2.05	0.57
3:C:2010:HEM:HBC2	3:C:2010:HEM:CMC	2.33	0.56
1:A:159:ILE:HD13	1:C:272:MET:HE1	1.86	0.56
3:B:1010:HEM:HBC2	3:B:1010:HEM:CMC	2.34	0.56
1:C:471:SER:O	1:C:477:HIS:HE1	1.87	0.56
1:D:100:LEU:CB	1:D:456:GLY:HA3	2.35	0.56
1:C:380:VAL:O	1:C:384:CYS:HB2	2.06	0.56
1:B:333:HIS:CD2	1:B:336:TYR:H	2.23	0.56
1:B:194:TRP:CZ3	1:B:206:TRP:HA	2.41	0.56
1:A:272:MET:HA	1:A:272:MET:HE2	1.88	0.56
1:D:130:ARG:NH1	1:D:133:PRO:HA	2.21	0.56
1:A:100:LEU:HD12	1:A:103:LYS:HD2	1.88	0.56
1:B:466:PRO:HG2	1:B:473:THR:HG21	1.88	0.56
1:D:352:VAL:CG2	5:D:3012:ITU:H11	2.35	0.55
1:C:194:TRP:CZ2	3:C:2010:HEM:HMC1	2.41	0.55
1:A:154:PHE:CD2	1:A:158:LYS:HD3	2.41	0.55
3:A:510:HEM:HHC	3:A:510:HEM:HBB2	1.89	0.55
1:B:354:ASN:HB2	7:B:1173:HOH:O	2.06	0.55
1:A:104:ALA:O	1:B:113:LYS:HD2	2.07	0.55
1:D:226:HIS:ND1	7:D:3344:HOH:O	2.32	0.55
1:B:105:LYS:HG2	1:B:484:VAL:O	2.08	0.54
1:B:357:LEU:HD22	1:B:359:VAL:HG22	1.88	0.54
1:B:387:GLN:HA	1:B:387:GLN:NE2	2.21	0.54
1:B:229:ARG:NH2	7:B:1031:HOH:O	2.40	0.54
1:A:104:ALA:HA	1:A:485:LEU:HD23	1.88	0.54
1:C:264:LEU:HD22	1:C:351:ALA:HB1	1.90	0.54
1:B:192:GLN:HB3	1:B:487:PRO:HG2	1.90	0.54
1:D:194:TRP:CZ2	3:D:3010:HEM:HBB2	2.43	0.54
7:A:3063:HOH:O	1:B:409:LEU:HA	2.07	0.54
1:C:194:TRP:CZ3	1:C:206:TRP:HA	2.43	0.54
1:A:196:ASN:O	1:A:198:PRO:HD3	2.08	0.54
1:A:374:MET:HE3	1:A:376:THR:OG1	2.08	0.54
1:B:226:HIS:HD2	7:B:1069:HOH:O	1.89	0.53
1:D:190:THR:HG22	1:D:357:LEU:HD11	1.88	0.53
1:A:263:GLN:NE2	1:A:266:ARG:HD3	2.22	0.53
1:B:194:TRP:CZ2	3:B:1010:HEM:HBB2	2.43	0.53
1:B:143:ALA:O	1:B:147:VAL:HG23	2.09	0.53
1:B:399:MET:HE3	1:B:417:GLU:HG3	1.90	0.53
1:C:381:ARG:NH2	6:C:2011:H4B:C4	2.71	0.53
1:D:156:GLU:HG2	1:D:158:LYS:HE2	1.90	0.53
1:A:194:TRP:CH2	3:A:510:HEM:HMC1	2.44	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:475:VAL:HG13	1:B:480:MET:HE1	1.90	0.53
1:A:369:PHE:CD1	5:A:512:ITU:H11	2.43	0.53
1:D:130:ARG:O	1:D:253:GLY:HA3	2.08	0.53
1:D:263:GLN:HE22	1:D:266:ARG:NE	2.07	0.53
1:B:232:ARG:HD2	7:B:1304:HOH:O	2.09	0.52
1:C:262:ALA:HA	1:C:354:ASN:ND2	2.24	0.52
1:B:226:HIS:HE1	7:B:1118:HOH:O	1.92	0.52
1:C:449:ASN:ND2	1:C:452:ARG:HH12	2.07	0.52
1:D:194:TRP:CH2	3:D:3010:HEM:HMC1	2.44	0.52
1:D:437:HIS:O	1:D:441:GLU:HG3	2.10	0.52
1:B:436:HIS:H	1:B:436:HIS:CD2	2.27	0.52
1:B:352:VAL:HB	1:B:369:PHE:CE1	2.45	0.52
1:C:205:GLN:HG3	3:C:2010:HEM:HBB1	1.90	0.52
1:D:374:MET:HE3	1:D:376:THR:OG1	2.09	0.52
1:B:440:ALA:HB1	1:B:475:VAL:HG23	1.92	0.52
1:D:381:ARG:O	1:D:385:ASP:HB2	2.10	0.52
1:A:167:GLU:O	1:A:171:LYS:HG2	2.09	0.52
1:A:193:ALA:HB2	1:A:487:PRO:HB2	1.91	0.52
1:A:204:ILE:HD11	1:A:445:LYS:HD3	1.92	0.52
1:A:83:ARG:O	1:A:102:HIS:CE1	2.62	0.52
1:A:86:ARG:NH2	1:A:95:THR:HG21	2.24	0.51
1:B:270:TYR:CE2	1:B:299:TYR:HA	2.46	0.51
3:D:3010:HEM:HBB2	3:D:3010:HEM:HHC	1.92	0.51
1:A:110:CYS:HB3	1:B:110:CYS:CB	2.39	0.51
1:B:263:GLN:HE21	1:B:266:ARG:HD3	1.76	0.51
1:B:475:VAL:CG1	1:B:480:MET:HE1	2.41	0.51
1:B:425:SER:O	1:B:429:GLN:HG2	2.11	0.51
1:D:475:VAL:HG13	1:D:480:MET:HE1	1.93	0.51
1:D:247:PHE:HB3	1:D:248:PRO:CD	2.41	0.50
1:B:194:TRP:HZ2	3:B:1010:HEM:HBB2	1.76	0.50
1:B:279:GLY:O	1:B:281:PRO:HD3	2.11	0.50
1:B:151:TYR:HA	1:B:154:PHE:CD2	2.46	0.50
1:B:162:HIS:O	1:B:166:VAL:HG23	2.12	0.50
1:B:271:GLN:HA	1:B:277:ILE:HD13	1.94	0.50
1:B:316:PRO:HD2	1:B:501:TRP:CZ3	2.47	0.50
1:B:357:LEU:HG	1:B:489:TYR:CE1	2.47	0.50
1:D:194:TRP:HZ2	3:D:3010:HEM:HBB2	1.76	0.50
1:A:394:GLU:HB2	7:A:3330:HOH:O	2.10	0.50
1:B:165:ARG:HA	7:B:1309:HOH:O	2.11	0.50
1:D:130:ARG:HH12	1:D:133:PRO:HA	1.77	0.50
1:C:156:GLU:H	1:C:156:GLU:CD	2.18	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:264:LEU:HB2	1:B:351:ALA:HB3	1.94	0.50
3:A:510:HEM:HBC2	3:A:510:HEM:CMC	2.42	0.50
1:B:333:HIS:NE2	1:B:335:LYS:HG2	2.27	0.50
1:C:451:TYR:CE1	1:C:456:GLY:HA2	2.46	0.50
1:D:449:ASN:HD22	1:D:452:ARG:HH12	1.59	0.49
1:C:88:LYS:HG3	1:C:95:THR:HG22	1.93	0.49
1:D:245:THR:O	1:D:367:CYS:HA	2.12	0.49
1:D:264:LEU:HD13	1:D:353:ALA:HB2	1.94	0.49
1:A:170:THR:O	1:A:174:GLU:HG3	2.12	0.49
1:A:194:TRP:CZ3	1:A:206:TRP:HA	2.47	0.49
1:B:377:GLU:HB3	7:B:1129:HOH:O	2.12	0.49
1:C:279:GLY:O	1:C:281:PRO:HD3	2.13	0.49
1:A:475:VAL:HG13	1:A:480:MET:HE1	1.95	0.49
1:B:192:GLN:HB3	1:B:487:PRO:CG	2.42	0.49
1:B:338:TRP:CE3	1:B:398:ARG:HD3	2.48	0.49
1:D:100:LEU:HB3	1:D:456:GLY:HA3	1.95	0.49
1:A:194:TRP:HZ2	3:A:510:HEM:HBB2	1.78	0.49
1:C:436:HIS:HD2	7:C:2061:HOH:O	1.94	0.49
1:B:151:TYR:CD1	1:B:154:PHE:CE2	3.00	0.49
1:C:412:ASP:O	1:C:416:VAL:HG23	2.13	0.49
1:D:350:PRO:O	1:D:350:PRO:HG2	2.13	0.49
1:A:186:LEU:HD22	1:A:215:ARG:HD2	1.95	0.49
3:A:510:HEM:HHC	3:A:510:HEM:CBB	2.43	0.49
1:C:352:VAL:HB	1:C:369:PHE:CE1	2.48	0.48
1:D:217:CYS:CB	7:D:3344:HOH:O	2.60	0.48
1:D:272:MET:HE1	1:D:299:TYR:HD1	1.78	0.48
1:D:338:TRP:CE3	1:D:398:ARG:HD3	2.48	0.48
1:A:393:GLU:HB2	7:A:3289:HOH:O	2.13	0.48
1:C:374:MET:HE3	1:C:376:THR:OG1	2.13	0.48
1:A:204:ILE:CD1	1:A:445:LYS:HD3	2.44	0.48
1:C:500:VAL:HA	7:C:2218:HOH:O	2.14	0.48
1:A:194:TRP:CZ2	3:A:510:HEM:HMC1	2.47	0.48
1:A:436:HIS:CD2	1:A:436:HIS:H	2.32	0.48
1:C:377:GLU:HB3	7:C:2131:HOH:O	2.14	0.48
1:D:304:VAL:HG21	1:D:326:VAL:HG11	1.95	0.48
1:A:333:HIS:CD2	1:A:336:TYR:H	2.31	0.48
1:D:107:ILE:HD11	1:D:484:VAL:HB	1.96	0.48
1:A:132:LYS:HG3	1:A:133:PRO:HD2	1.95	0.48
1:A:416:VAL:HG11	1:B:413:GLN:HG2	1.96	0.48
1:A:423:ILE:O	1:A:427:GLN:HG3	2.13	0.48
1:C:111:ARG:HH21	1:D:101:HIS:CD2	2.31	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:399:MET:CE	1:D:417:GLU:HG3	2.44	0.48
1:B:182:THR:OG1	1:B:185:GLU:HG3	2.13	0.48
1:B:356:LEU:C	1:B:356:LEU:HD23	2.39	0.48
1:B:416:VAL:O	1:B:420:ILE:HG13	2.13	0.48
1:A:194:TRP:CZ2	3:A:510:HEM:HBB2	2.49	0.47
1:C:248:PRO:HD2	1:C:257:PHE:CG	2.49	0.47
1:D:411:LYS:O	1:D:415:VAL:HG12	2.14	0.47
1:C:100:LEU:HD12	1:C:103:LYS:HD2	1.97	0.47
1:D:310:GLN:HG3	1:D:314:ARG:O	2.14	0.47
1:A:110:CYS:CB	1:B:110:CYS:HB3	2.39	0.47
1:B:270:TYR:CD2	1:B:299:TYR:HA	2.50	0.47
1:B:435:ASP:OD2	1:B:437:HIS:HB2	2.14	0.47
1:A:263:GLN:HE22	1:A:266:ARG:NH1	2.13	0.47
1:C:298:LYS:HE2	7:C:2100:HOH:O	2.14	0.47
1:A:230:HIS:ND1	1:A:245:THR:OG1	2.42	0.47
7:A:3017:HOH:O	1:B:477:HIS:HD2	1.96	0.47
1:B:199:ARG:HD3	1:B:463:TRP:CD2	2.49	0.47
1:B:337:GLU:O	1:B:340:ARG:HG2	2.15	0.47
1:B:316:PRO:HD2	1:B:501:TRP:CH2	2.49	0.47
1:C:100:LEU:HB3	1:C:456:GLY:HA3	1.95	0.47
1:C:131:ASP:OD1	1:C:132:LYS:HG3	2.15	0.47
1:C:449:ASN:HD22	1:C:452:ARG:HH12	1.61	0.47
1:D:295:TRP:HB2	1:D:318:LEU:HD23	1.95	0.47
3:D:3010:HEM:HHC	3:D:3010:HEM:CBB	2.44	0.47
1:A:322:PRO:HD2	1:A:325:LEU:HD12	1.97	0.47
1:C:130:ARG:NH1	1:C:133:PRO:HA	2.30	0.47
1:C:463:TRP:HZ3	3:C:2010:HEM:HBA2	1.79	0.47
1:A:107:ILE:HD11	1:A:484:VAL:HG11	1.97	0.47
1:A:156:GLU:O	1:A:156:GLU:HG2	2.14	0.47
1:A:333:HIS:HD2	1:A:335:LYS:H	1.63	0.47
1:A:423:ILE:HD11	1:B:409:LEU:HG	1.97	0.47
1:A:247:PHE:HB3	1:A:248:PRO:CD	2.44	0.47
1:A:338:TRP:CE3	1:A:398:ARG:HD3	2.50	0.47
1:B:335:LYS:HE3	1:B:336:TYR:CE1	2.50	0.47
1:C:156:GLU:CD	1:C:156:GLU:N	2.73	0.46
1:C:236:ASN:O	1:C:237:ASN:HB2	2.15	0.46
1:A:140:LEU:O	1:A:144:ILE:HG12	2.15	0.46
1:B:149:GLN:NE2	7:B:1122:HOH:O	2.48	0.46
1:B:394:GLU:HG3	7:B:1337:HOH:O	2.16	0.46
1:C:245:THR:O	1:C:367:CYS:HA	2.15	0.46
1:C:130:ARG:HD3	7:C:2269:HOH:O	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:191:LYS:O	1:C:194:TRP:HB3	2.16	0.46
1:A:88:LYS:CB	1:A:95:THR:HG22	2.39	0.46
1:D:272:MET:HE1	1:D:299:TYR:CD1	2.51	0.46
1:B:463:TRP:HA	6:B:1011:H4B:N1	2.31	0.46
1:A:380:VAL:O	1:A:384:CYS:HB2	2.16	0.46
1:B:391:ILE:O	1:B:395:VAL:HG23	2.15	0.46
1:A:263:GLN:HG2	7:A:3122:HOH:O	2.15	0.46
1:D:159:ILE:HG22	1:D:163:LEU:HD12	1.98	0.46
1:A:333:HIS:CD2	1:A:335:LYS:H	2.34	0.46
1:A:471:SER:O	1:A:477:HIS:CE1	2.64	0.46
1:B:140:LEU:O	1:B:144:ILE:HG12	2.15	0.46
1:C:247:PHE:HB3	1:C:248:PRO:CD	2.45	0.46
1:A:85:VAL:HG11	1:A:481:LEU:HD21	1.98	0.45
1:C:101:HIS:O	1:D:113:LYS:NZ	2.49	0.45
1:C:255:HIS:HA	1:C:313:GLY:HA3	1.98	0.45
1:C:467:PRO:HB2	1:D:471:SER:OG	2.15	0.45
3:D:3010:HEM:HBC2	3:D:3010:HEM:CMC	2.37	0.45
1:A:269:GLY:O	1:A:284:VAL:HA	2.16	0.45
1:A:446:TYR:CZ	1:A:450:GLU:HG3	2.50	0.45
1:C:322:PRO:HA	1:C:323:PRO:HD3	1.75	0.45
1:D:226:HIS:CE1	7:D:3344:HOH:O	2.70	0.45
1:D:332:GLU:O	1:D:424:HIS:HD2	1.99	0.45
1:C:194:TRP:CE2	3:C:2010:HEM:C2C	3.05	0.45
1:A:204:ILE:HG13	1:A:446:TYR:HB2	1.97	0.45
1:D:463:TRP:HA	6:D:3011:H4B:N1	2.32	0.45
1:A:158:LYS:HB3	1:A:158:LYS:NZ	2.31	0.45
1:B:308:VAL:HG12	1:B:316:PRO:HB2	1.99	0.45
1:C:193:ALA:HB2	1:C:487:PRO:HB2	1.99	0.45
1:C:338:TRP:CZ3	1:C:398:ARG:HD3	2.52	0.45
1:C:394:GLU:HB2	7:C:2048:HOH:O	2.16	0.45
1:B:85:VAL:HB	1:B:481:LEU:HD11	1.98	0.45
1:B:151:TYR:HA	1:B:154:PHE:CE2	2.51	0.45
1:C:335:LYS:HE2	7:C:2240:HOH:O	2.16	0.45
1:A:375:GLY:HA2	1:A:433:ILE:HD11	1.99	0.45
1:B:263:GLN:HE22	1:B:266:ARG:NH1	2.15	0.45
1:B:449:ASN:HD22	1:B:452:ARG:NH1	2.15	0.45
1:B:451:TYR:CE1	1:B:456:GLY:HA2	2.52	0.45
1:C:155:LYS:HB2	1:C:156:GLU:OE1	2.17	0.45
1:D:107:ILE:HD11	1:D:484:VAL:CB	2.47	0.45
1:B:137:ASP:O	1:B:141:PRO:HD2	2.17	0.44
1:A:183:GLY:O	1:A:187:ILE:HG13	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:116:LEU:N	1:B:116:LEU:HD23	2.32	0.44
1:D:151:TYR:CD1	1:D:154:PHE:CE2	3.06	0.44
1:D:217:CYS:HB2	7:D:3344:HOH:O	2.16	0.44
1:A:116:LEU:O	1:A:119:ILE:HG12	2.17	0.44
1:C:230:HIS:C	1:C:230:HIS:CD2	2.95	0.44
1:C:423:ILE:O	1:C:427:GLN:HG3	2.18	0.44
1:A:190:THR:HG22	1:A:357:LEU:HD11	2.00	0.44
1:B:245:THR:O	1:B:367:CYS:HA	2.18	0.44
1:D:412:ASP:O	1:D:416:VAL:HG23	2.18	0.44
7:C:2056:HOH:O	1:D:477:HIS:HD2	2.00	0.44
1:A:502:GLN:HG2	1:A:502:GLN:O	2.18	0.44
1:B:471:SER:O	1:B:477:HIS:HE1	2.00	0.44
1:C:159:ILE:H	1:C:159:ILE:CD1	2.27	0.44
1:C:374:MET:HA	1:C:434:MET:O	2.18	0.44
1:B:196:ASN:O	1:B:198:PRO:HD3	2.18	0.44
1:D:338:TRP:O	1:D:341:GLU:HB2	2.18	0.44
1:A:279:GLY:O	1:A:281:PRO:HD3	2.18	0.43
1:B:232:ARG:CZ	1:B:325:LEU:HD21	2.48	0.43
1:C:260:TRP:HB2	1:C:308:VAL:HB	1.99	0.43
1:C:405:LYS:HE2	1:C:405:LYS:HB3	1.78	0.43
1:A:111:ARG:HH12	1:A:113:LYS:HB3	1.80	0.43
1:B:263:GLN:NE2	1:B:266:ARG:HD3	2.32	0.43
1:C:135:PRO:HA	1:C:136:PRO:HD3	1.81	0.43
1:C:159:ILE:HD13	1:C:159:ILE:N	2.27	0.43
1:D:348:ALA:HB1	1:D:431:VAL:HG11	2.00	0.43
1:A:116:LEU:N	1:A:116:LEU:HD23	2.33	0.43
1:A:444:MET:HE1	1:A:478:GLN:OE1	2.18	0.43
1:B:448:GLN:HG2	1:B:452:ARG:NH2	2.33	0.43
1:D:107:ILE:HD11	1:D:484:VAL:HG11	2.00	0.43
1:D:194:TRP:CZ2	3:D:3010:HEM:HMC1	2.53	0.43
1:B:130:ARG:HB2	1:B:250:ARG:HD3	1.99	0.43
1:C:321:ILE:HG13	7:C:2026:HOH:O	2.18	0.43
1:C:499:HIS:HD2	1:C:501:TRP:CE2	2.36	0.43
1:B:291:ILE:HD11	1:B:297:PRO:HB3	2.00	0.43
1:C:87:ILE:HG13	1:C:96:PHE:HB2	1.99	0.43
1:C:295:TRP:HB2	1:C:318:LEU:HD23	2.00	0.43
1:C:387:GLN:HB3	7:C:2259:HOH:O	2.18	0.43
1:D:357:LEU:HG	1:D:489:TYR:CE1	2.54	0.43
1:B:130:ARG:HD2	1:B:132:LYS:O	2.18	0.43
1:B:357:LEU:HD22	1:B:359:VAL:CG2	2.47	0.43
1:B:380:VAL:O	1:B:384:CYS:HB2	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:427:GLN:HA	7:D:3089:HOH:O	2.18	0.43
1:B:306:PRO:HB3	1:B:320:GLU:HG2	2.00	0.43
1:D:358:GLU:HA	1:D:362:LEU:O	2.19	0.43
1:C:475:VAL:HG13	1:C:480:MET:HE1	1.99	0.43
3:C:2010:HEM:C4C	5:C:2012:ITU:H13	2.54	0.43
1:B:270:TYR:O	1:B:277:ILE:HA	2.18	0.42
1:B:278:ARG:NH1	1:B:301:ARG:HG3	2.33	0.42
1:C:349:LEU:HD11	1:C:370:ASN:ND2	2.33	0.42
1:A:111:ARG:HH11	1:A:111:ARG:CB	2.26	0.42
1:A:451:TYR:CE1	1:A:456:GLY:HA2	2.54	0.42
1:D:84:HIS:CE1	1:D:97:GLN:HG2	2.54	0.42
1:D:357:LEU:HB3	1:D:364:PHE:HB2	2.00	0.42
1:A:291:ILE:CD1	1:A:297:PRO:HB3	2.49	0.42
1:D:127:ARG:NH1	1:D:358:GLU:HG3	2.34	0.42
1:C:116:LEU:HD12	7:C:2222:HOH:O	2.19	0.42
1:D:266:ARG:HD3	7:D:3040:HOH:O	2.18	0.42
1:A:186:LEU:HD11	1:A:246:VAL:HG21	2.01	0.42
1:A:232:ARG:HD3	7:A:3211:HOH:O	2.19	0.42
1:A:349:LEU:HD11	1:A:370:ASN:ND2	2.35	0.42
1:B:369:PHE:CD1	5:B:1012:ITU:C1	3.02	0.42
1:D:190:THR:HG22	1:D:357:LEU:CD1	2.49	0.42
1:D:399:MET:HE1	1:D:417:GLU:HG3	2.02	0.42
1:B:233:TYR:O	1:B:236:ASN:HB2	2.19	0.42
1:B:247:PHE:HB3	1:B:248:PRO:CD	2.50	0.42
1:C:169:VAL:O	1:C:173:ILE:HG13	2.19	0.42
1:C:425:SER:O	1:C:429:GLN:HG2	2.19	0.42
1:C:107:ILE:HD12	1:C:124:SER:O	2.19	0.42
1:D:258:ARG:HG2	1:D:365:PRO:HB2	2.01	0.42
1:C:301:ARG:HH22	1:C:343:GLU:HB3	1.84	0.42
1:D:264:LEU:HD12	1:D:264:LEU:HA	1.90	0.42
1:A:199:ARG:HD3	1:A:463:TRP:CE2	2.55	0.42
1:B:369:PHE:HD1	5:B:1012:ITU:C1	2.33	0.42
1:A:270:TYR:CE2	1:A:299:TYR:HA	2.55	0.41
1:B:181:LEU:O	1:B:215:ARG:NH2	2.47	0.41
1:B:263:GLN:HA	1:B:351:ALA:O	2.20	0.41
1:D:265:ILE:HD12	1:D:349:LEU:HB3	2.02	0.41
1:A:459:ALA:HB1	1:A:464:LEU:CD1	2.50	0.41
1:B:203:ARG:O	1:B:206:TRP:HD1	2.02	0.41
1:B:263:GLN:HG2	7:B:1130:HOH:O	2.19	0.41
1:C:180:GLN:NE2	7:C:2210:HOH:O	2.53	0.41
1:C:216:SER:HB2	7:C:2311:HOH:O	2.18	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:233:TYR:CD1	1:C:233:TYR:C	2.98	0.41
1:C:308:VAL:HG12	1:C:316:PRO:HB2	2.02	0.41
1:C:383:PHE:O	1:C:391:ILE:HG12	2.20	0.41
1:D:475:VAL:HG13	1:D:480:MET:CE	2.51	0.41
3:A:510:HEM:HBB2	3:A:510:HEM:CHC	2.50	0.41
1:A:149:GLN:HG2	1:A:188:PHE:HZ	1.84	0.41
1:A:392:LEU:HB2	7:A:3092:HOH:O	2.20	0.41
1:B:160:GLU:CD	1:B:160:GLU:H	2.28	0.41
1:B:449:ASN:ND2	1:B:452:ARG:NH1	2.67	0.41
1:C:116:LEU:HB3	1:D:481:LEU:CD2	2.51	0.41
1:C:301:ARG:HH22	1:C:343:GLU:CB	2.33	0.41
1:B:322:PRO:HA	1:B:323:PRO:HD2	1.69	0.41
1:B:418:ILE:O	1:B:421:ALA:HB3	2.20	0.41
1:B:465:VAL:HA	1:B:466:PRO:HD3	1.96	0.41
1:D:194:TRP:CE2	3:D:3010:HEM:C2C	3.08	0.41
1:A:158:LYS:HB2	7:A:3115:HOH:O	2.20	0.41
1:C:197:ALA:HA	1:C:198:PRO:HD3	1.90	0.41
1:C:383:PHE:HA	1:C:389:TYR:HB2	2.03	0.41
1:A:352:VAL:HB	1:A:369:PHE:CE1	2.55	0.41
1:C:173:ILE:O	1:C:177:GLY:N	2.52	0.41
1:C:316:PRO:HD2	1:C:501:TRP:CH2	2.56	0.41
1:C:336:TYR:HB3	1:C:338:TRP:CE2	2.56	0.41
1:D:381:ARG:NH1	1:D:385:ASP:OD2	2.53	0.41
1:B:104:ALA:CB	1:B:483:TYR:HB2	2.51	0.41
1:C:85:VAL:HB	1:C:481:LEU:HD11	2.03	0.41
1:C:161:GLU:HG3	7:C:2128:HOH:O	2.20	0.41
1:D:83:ARG:HH21	1:D:84:HIS:HB3	1.86	0.41
1:D:263:GLN:NE2	1:D:266:ARG:HG2	2.36	0.41
1:D:471:SER:HA	1:D:476:PHE:CG	2.56	0.41
1:A:332:GLU:HG2	1:A:340:ARG:HD3	2.02	0.40
1:C:309:LEU:O	1:C:316:PRO:HA	2.21	0.40
1:C:481:LEU:CD2	1:D:116:LEU:HD22	2.38	0.40
1:D:279:GLY:HA2	1:D:301:ARG:HA	2.01	0.40
1:D:296:LYS:HA	1:D:297:PRO:HD2	1.90	0.40
1:A:377:GLU:O	1:A:381:ARG:HB2	2.20	0.40
1:C:306:PRO:HB3	1:C:320:GLU:HG2	2.04	0.40
1:D:249:GLN:HB3	1:D:364:PHE:CE2	2.56	0.40
1:D:333:HIS:CD2	1:D:334:PRO:HD2	2.56	0.40
1:A:248:PRO:HD2	1:A:257:PHE:CG	2.57	0.40
1:A:293:LEU:HD21	1:A:501:TRP:CH2	2.57	0.40
1:C:196:ASN:O	1:C:198:PRO:HD3	2.20	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:392:LEU:HB2	7:C:2104:HOH:O	2.21	0.40
1:B:171:LYS:O	1:B:174:GLU:HB3	2.21	0.40
1:C:273:PRO:C	1:C:275:GLY:H	2.28	0.40
1:C:416:VAL:HG11	1:D:413:GLN:HG2	2.04	0.40
1:C:436:HIS:H	1:C:436:HIS:CD2	2.39	0.40
1:B:115:CYS:SG	7:B:1220:HOH:O	2.63	0.40
1:D:333:HIS:CD2	1:D:335:LYS:HG3	2.57	0.40

All (1) 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)
1:C:137:ASP:O	7:B:1165:HOH:O[4_455]	2.19	0.01

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	419/427 (98%)	390 (93%)	28 (7%)	1 (0%)	43	50
1	B	419/427 (98%)	395 (94%)	24 (6%)	0	100	100
1	C	419/427 (98%)	390 (93%)	26 (6%)	3 (1%)	18	17
1	D	419/427 (98%)	398 (95%)	19 (4%)	2 (0%)	24	24
All	All	1676/1708 (98%)	1573 (94%)	97 (6%)	6 (0%)	30	31

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	236	ASN
1	D	206	TRP
1	C	206	TRP

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Mol	Chain	Res	Type
1	A	203	ARG
1	D	236	ASN
1	C	275	GLY

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	366/373 (98%)	350 (96%)	16 (4%)	25	29
1	B	366/373 (98%)	347 (95%)	19 (5%)	21	22
1	C	366/373 (98%)	346 (94%)	20 (6%)	19	20
1	D	366/373 (98%)	348 (95%)	18 (5%)	22	24
All	All	1464/1492 (98%)	1391 (95%)	73 (5%)	22	24

All (73) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	88	LYS
1	A	111	ARG
1	A	116	LEU
1	A	158	LYS
1	A	167	GLU
1	A	181	LEU
1	A	264	LEU
1	A	296	LYS
1	A	323	PRO
1	A	337	GLU
1	A	340	ARG
1	A	357	LEU
1	A	362	LEU
1	A	393	GLU
1	A	423	ILE
1	A	502	GLN
1	B	113	LYS
1	B	116	LEU

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Mol	Chain	Res	Type
1	B	155	LYS
1	B	156	GLU
1	B	163	LEU
1	B	181	LEU
1	B	216	SER
1	B	217	CYS
1	B	264	LEU
1	B	285	GLU
1	B	337	GLU
1	B	357	LEU
1	B	362	LEU
1	B	387	GLN
1	B	393	GLU
1	B	423	ILE
1	B	464	LEU
1	B	494	GLU
1	B	502	GLN
1	C	116	LEU
1	C	145	GLU
1	C	155	LYS
1	C	156	GLU
1	C	159	ILE
1	C	167	GLU
1	C	180	GLN
1	C	181	LEU
1	C	264	LEU
1	C	274	ASP
1	C	278	ARG
1	C	285	GLU
1	C	324	ASP
1	C	337	GLU
1	C	362	LEU
1	C	381	ARG
1	C	393	GLU
1	C	394	GLU
1	C	423	ILE
1	C	464	LEU
1	D	97	GLN
1	D	116	LEU
1	D	123	LYS
1	D	137	ASP
1	D	160	GLU

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Mol	Chain	Res	Type
1	D	171	LYS
1	D	174	GLU
1	D	181	LEU
1	D	221	ARG
1	D	242	SER
1	D	263	GLN
1	D	264	LEU
1	D	266	ARG
1	D	278	ARG
1	D	320	GLU
1	D	357	LEU
1	D	362	LEU
1	D	464	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (60) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	180	GLN
1	A	205	GLN
1	A	208	ASN
1	A	226	HIS
1	A	312	ASN
1	A	333	HIS
1	A	370	ASN
1	A	390	ASN
1	A	404	HIS
1	A	413	GLN
1	A	424	HIS
1	A	436	HIS
1	A	448	GLN
1	A	449	ASN
1	A	477	HIS
1	B	102	HIS
1	B	149	GLN
1	B	180	GLN
1	B	205	GLN
1	B	208	ASN
1	B	226	HIS
1	B	239	ASN
1	B	263	GLN
1	B	271	GLN
1	B	333	HIS

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Mol	Chain	Res	Type
1	B	370	ASN
1	B	387	GLN
1	B	390	ASN
1	B	413	GLN
1	B	436	HIS
1	B	448	GLN
1	B	449	ASN
1	B	477	HIS
1	B	492	GLN
1	B	502	GLN
1	C	102	HIS
1	C	149	GLN
1	C	180	GLN
1	C	226	HIS
1	C	271	GLN
1	C	333	HIS
1	C	436	HIS
1	C	449	ASN
1	C	477	HIS
1	C	492	GLN
1	C	502	GLN
1	D	102	HIS
1	D	192	GLN
1	D	205	GLN
1	D	208	ASN
1	D	226	HIS
1	D	263	GLN
1	D	310	GLN
1	D	333	HIS
1	D	413	GLN
1	D	429	GLN
1	D	448	GLN
1	D	449	ASN
1	D	477	HIS
1	D	492	GLN

5.3.3 RNA ⓘ

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](#)

Of 14 ligands modelled in this entry, 2 are monoatomic - leaving 12 for Mogul analysis.

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
3	HEM	D	3010	1	50,50,50	1.84	10 (20%)	67,82,82	1.01	4 (5%)
4	H2B	A	511	-	15,18,18	3.49	4 (26%)	15,26,26	2.14	4 (26%)
5	ITU	B	1012	-	4,5,5	3.48	1 (25%)	3,5,5	1.83	1 (33%)
6	H4B	B	1011	-	17,18,18	1.86	3 (17%)	14,26,26	1.75	3 (21%)
6	H4B	C	2011	-	17,18,18	1.55	3 (17%)	14,26,26	1.86	4 (28%)
6	H4B	D	3011	-	17,18,18	1.80	4 (23%)	14,26,26	1.82	3 (21%)
5	ITU	D	3012	-	4,5,5	3.68	1 (25%)	3,5,5	0.75	0
5	ITU	C	2012	-	4,5,5	2.92	1 (25%)	3,5,5	1.71	1 (33%)
3	HEM	B	1010	1	50,50,50	1.79	11 (22%)	67,82,82	1.11	5 (7%)
3	HEM	A	510	1	50,50,50	2.00	9 (18%)	67,82,82	1.08	6 (8%)
5	ITU	A	512	-	4,5,5	2.89	1 (25%)	3,5,5	1.39	1 (33%)
3	HEM	C	2010	1	50,50,50	2.02	9 (18%)	67,82,82	1.12	5 (7%)

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
3	HEM	D	3010	1	-	3/14/54/54	-
4	H2B	A	511	-	-	0/8/33/33	0/2/2/2
5	ITU	B	1012	-	-	0/3/3/3	-
6	H4B	B	1011	-	-	0/8/17/17	0/2/2/2
6	H4B	C	2011	-	-	0/8/17/17	0/2/2/2
6	H4B	D	3011	-	-	0/8/17/17	0/2/2/2
5	ITU	D	3012	-	-	0/3/3/3	-
5	ITU	C	2012	-	-	0/3/3/3	-
3	HEM	B	1010	1	-	2/14/54/54	-
3	HEM	A	510	1	-	5/14/54/54	-
5	ITU	A	512	-	-	0/3/3/3	-
3	HEM	C	2010	1	-	6/14/54/54	-

All (57) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	511	H2B	C4A-N5	11.84	1.39	1.28
5	D	3012	ITU	C3-S	7.19	1.83	1.75
3	C	2010	HEM	FE-NB	6.87	2.16	1.94
5	B	1012	ITU	C3-S	6.84	1.82	1.75
5	A	512	ITU	C3-S	5.68	1.81	1.75
5	C	2012	ITU	C3-S	5.61	1.81	1.75
3	B	1010	HEM	FE-NC	5.11	2.12	1.95
3	D	3010	HEM	CBC-CAC	4.98	1.54	1.30
3	A	510	HEM	FE-NA	4.96	2.11	1.95
3	D	3010	HEM	FE-NA	4.96	2.11	1.95
3	A	510	HEM	FE-ND	4.93	2.10	1.94
3	A	510	HEM	CBB-CAB	4.90	1.53	1.30
6	B	1011	H4B	C7-C6	-4.85	1.47	1.52
3	C	2010	HEM	FE-ND	4.81	2.09	1.94
3	B	1010	HEM	CBB-CAB	4.80	1.53	1.30
3	C	2010	HEM	CBB-CAB	4.79	1.53	1.30
3	D	3010	HEM	CBB-CAB	4.78	1.53	1.30
3	A	510	HEM	CBC-CAC	4.77	1.53	1.30
3	B	1010	HEM	CBC-CAC	4.74	1.53	1.30
3	C	2010	HEM	CBC-CAC	4.66	1.52	1.30
3	A	510	HEM	FE-NC	4.63	2.10	1.95
3	D	3010	HEM	FE-NB	4.57	2.09	1.94
6	D	3011	H4B	C4A-C8A	4.54	1.46	1.38
3	A	510	HEM	FE-NB	4.43	2.08	1.94
6	B	1011	H4B	C4A-C8A	3.95	1.45	1.38
3	C	2010	HEM	FE-NA	3.91	2.08	1.95

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	C	2011	H4B	C4A-C8A	3.64	1.44	1.38
4	A	511	H2B	C8A-N1	3.58	1.39	1.32
3	B	1010	HEM	FE-NA	3.57	2.07	1.95
3	C	2010	HEM	FE-NC	3.52	2.06	1.95
3	D	3010	HEM	FE-NC	3.50	2.06	1.95
6	D	3011	H4B	C7-C6	-3.48	1.48	1.52
6	C	2011	H4B	C7-C6	-3.48	1.48	1.52
4	A	511	H2B	C4-N3	3.41	1.43	1.38
3	D	3010	HEM	CAC-C3C	3.04	1.55	1.47
3	C	2010	HEM	CAB-C3B	2.89	1.55	1.47
3	B	1010	HEM	CAC-C3C	2.88	1.55	1.47
3	A	510	HEM	CAB-C3B	2.85	1.55	1.47
3	A	510	HEM	CAC-C3C	2.80	1.54	1.47
3	B	1010	HEM	CAB-C3B	2.78	1.54	1.47
6	D	3011	H4B	C6-N5	2.75	1.52	1.47
3	C	2010	HEM	O2A-CGA	-2.73	1.21	1.30
3	A	510	HEM	O2A-CGA	-2.68	1.21	1.30
3	C	2010	HEM	CAC-C3C	2.62	1.54	1.47
3	D	3010	HEM	O2A-CGA	-2.52	1.22	1.30
3	B	1010	HEM	FE-NB	2.44	2.02	1.94
3	B	1010	HEM	O1A-CGA	2.35	1.29	1.22
3	B	1010	HEM	FE-ND	2.33	2.02	1.94
3	D	3010	HEM	O1A-CGA	2.20	1.29	1.22
4	A	511	H2B	C4-C4A	-2.15	1.41	1.47
6	C	2011	H4B	C6-N5	2.15	1.51	1.47
3	D	3010	HEM	CAB-C3B	2.14	1.53	1.47
6	B	1011	H4B	C2-N1	2.14	1.38	1.33
6	D	3011	H4B	C4-N3	2.14	1.42	1.38
3	B	1010	HEM	O2A-CGA	-2.06	1.24	1.30
3	B	1010	HEM	CMD-C2D	2.04	1.54	1.50
3	D	3010	HEM	CMC-C2C	2.01	1.54	1.50

All (37) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	C	2011	H4B	C2-N1-C8A	5.43	122.95	113.36
6	D	3011	H4B	C2-N1-C8A	5.26	122.64	113.36
4	A	511	H2B	C4-C4A-C8A	5.12	123.21	117.40
6	B	1011	H4B	C2-N1-C8A	5.08	122.33	113.36
4	A	511	H2B	N1-C2-N3	-4.03	120.05	126.44
4	A	511	H2B	C4-C4A-N5	3.32	121.65	118.06
3	C	2010	HEM	CBD-CAD-C3D	-3.14	103.85	112.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1010	HEM	CBD-CAD-C3D	-3.13	103.88	112.53
5	B	1012	ITU	C2-S-C3	-2.99	100.04	103.15
3	B	1010	HEM	O1A-CGA-CBA	-2.73	114.44	123.09
5	C	2012	ITU	C2-S-C3	-2.71	100.33	103.15
3	A	510	HEM	CBD-CAD-C3D	-2.71	105.04	112.53
3	A	510	HEM	O1A-CGA-CBA	-2.71	114.51	123.09
3	D	3010	HEM	CBB-CAB-C3B	-2.53	114.87	127.53
3	D	3010	HEM	CBD-CAD-C3D	-2.50	105.62	112.53
3	C	2010	HEM	O1A-CGA-CBA	-2.47	115.25	123.09
3	A	510	HEM	CBB-CAB-C3B	-2.43	115.37	127.53
3	B	1010	HEM	CBC-CAC-C3C	-2.42	115.41	127.53
3	D	3010	HEM	O1A-CGA-CBA	-2.38	115.53	123.09
3	C	2010	HEM	CBC-CAC-C3C	-2.38	115.62	127.53
3	B	1010	HEM	CBB-CAB-C3B	-2.32	115.95	127.53
3	D	3010	HEM	CBC-CAC-C3C	-2.27	116.18	127.53
6	D	3011	H4B	C4-C4A-N5	2.23	122.34	116.27
5	A	512	ITU	C2-S-C3	-2.22	100.83	103.15
3	A	510	HEM	C2A-C1A-NA	-2.22	107.69	110.15
6	C	2011	H4B	N2-C2-N3	2.21	121.42	116.76
3	C	2010	HEM	CBB-CAB-C3B	-2.15	116.77	127.53
3	B	1010	HEM	CHB-C1B-NB	2.15	127.02	124.37
3	A	510	HEM	CBC-CAC-C3C	-2.14	116.81	127.53
6	D	3011	H4B	N2-C2-N3	2.13	121.25	116.76
6	C	2011	H4B	N3-C2-N1	-2.12	119.44	123.32
4	A	511	H2B	N2-C2-N3	2.09	120.32	117.09
6	B	1011	H4B	N3-C2-N1	-2.04	119.59	123.32
6	C	2011	H4B	C4-C4A-N5	2.04	121.83	116.27
6	B	1011	H4B	C4-C4A-N5	2.04	121.83	116.27
3	C	2010	HEM	C2A-C1A-NA	-2.02	107.91	110.15
3	A	510	HEM	CBA-CAA-C2A	-2.00	106.99	112.53

There are no chirality outliers.

All (16) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	1010	HEM	C2A-CAA-CBA-CGA
3	D	3010	HEM	C2A-CAA-CBA-CGA
3	C	2010	HEM	C2A-CAA-CBA-CGA
3	B	1010	HEM	C1A-C2A-CAA-CBA
3	C	2010	HEM	CAA-CBA-CGA-O2A
3	A	510	HEM	C2A-CAA-CBA-CGA
3	C	2010	HEM	C1A-C2A-CAA-CBA

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Mol	Chain	Res	Type	Atoms
3	C	2010	HEM	CAA-CBA-CGA-O1A
3	A	510	HEM	CAA-CBA-CGA-O1A
3	A	510	HEM	CAA-CBA-CGA-O2A
3	D	3010	HEM	CAD-CBD-CGD-O2D
3	C	2010	HEM	CAD-CBD-CGD-O2D
3	A	510	HEM	CAD-CBD-CGD-O2D
3	C	2010	HEM	C2B-C3B-CAB-CBB
3	D	3010	HEM	CAA-CBA-CGA-O2A
3	A	510	HEM	CAD-CBD-CGD-O1D

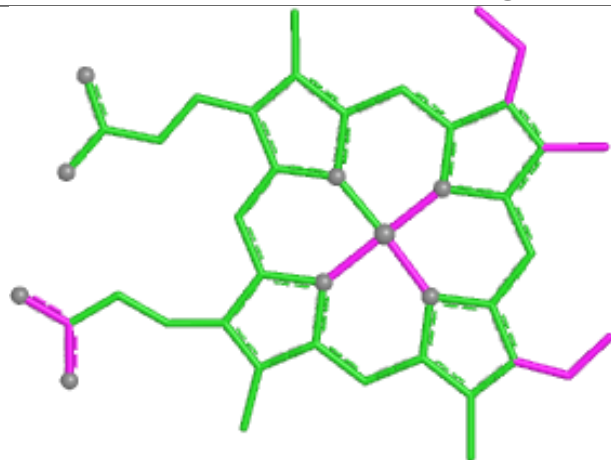
There are no ring outliers.

11 monomers are involved in 44 short contacts:

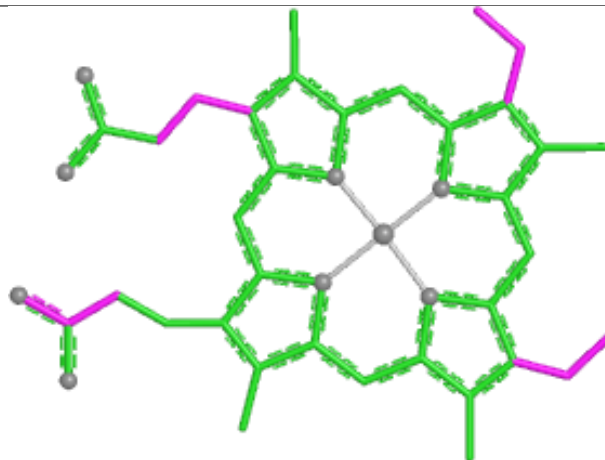
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	3010	HEM	9	0
5	B	1012	ITU	4	0
6	B	1011	H4B	1	0
6	C	2011	H4B	2	0
6	D	3011	H4B	1	0
5	D	3012	ITU	2	0
5	C	2012	ITU	3	0
3	B	1010	HEM	4	0
3	A	510	HEM	9	0
5	A	512	ITU	2	0
3	C	2010	HEM	8	0

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.

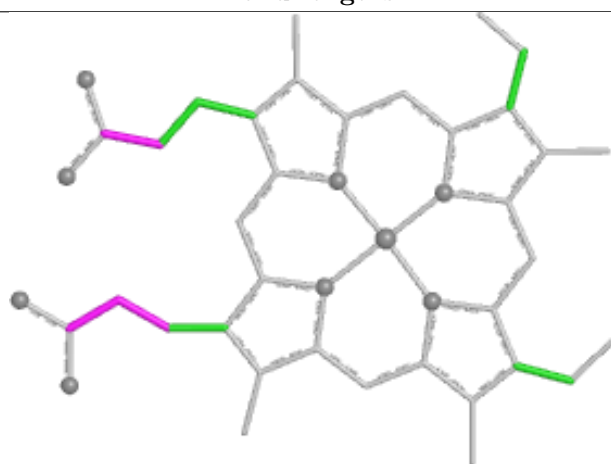
Ligand HEM D 3010



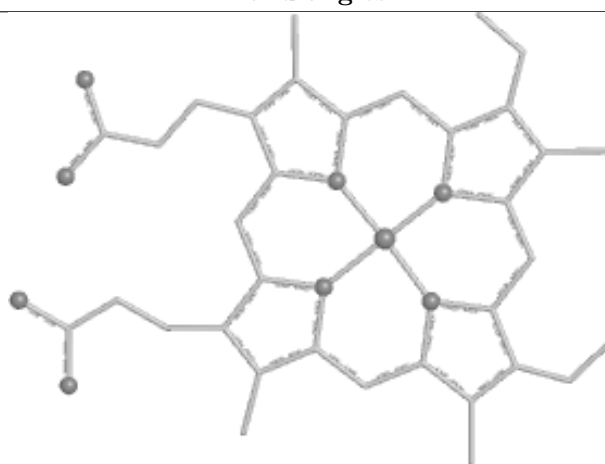
Bond lengths



Bond angles

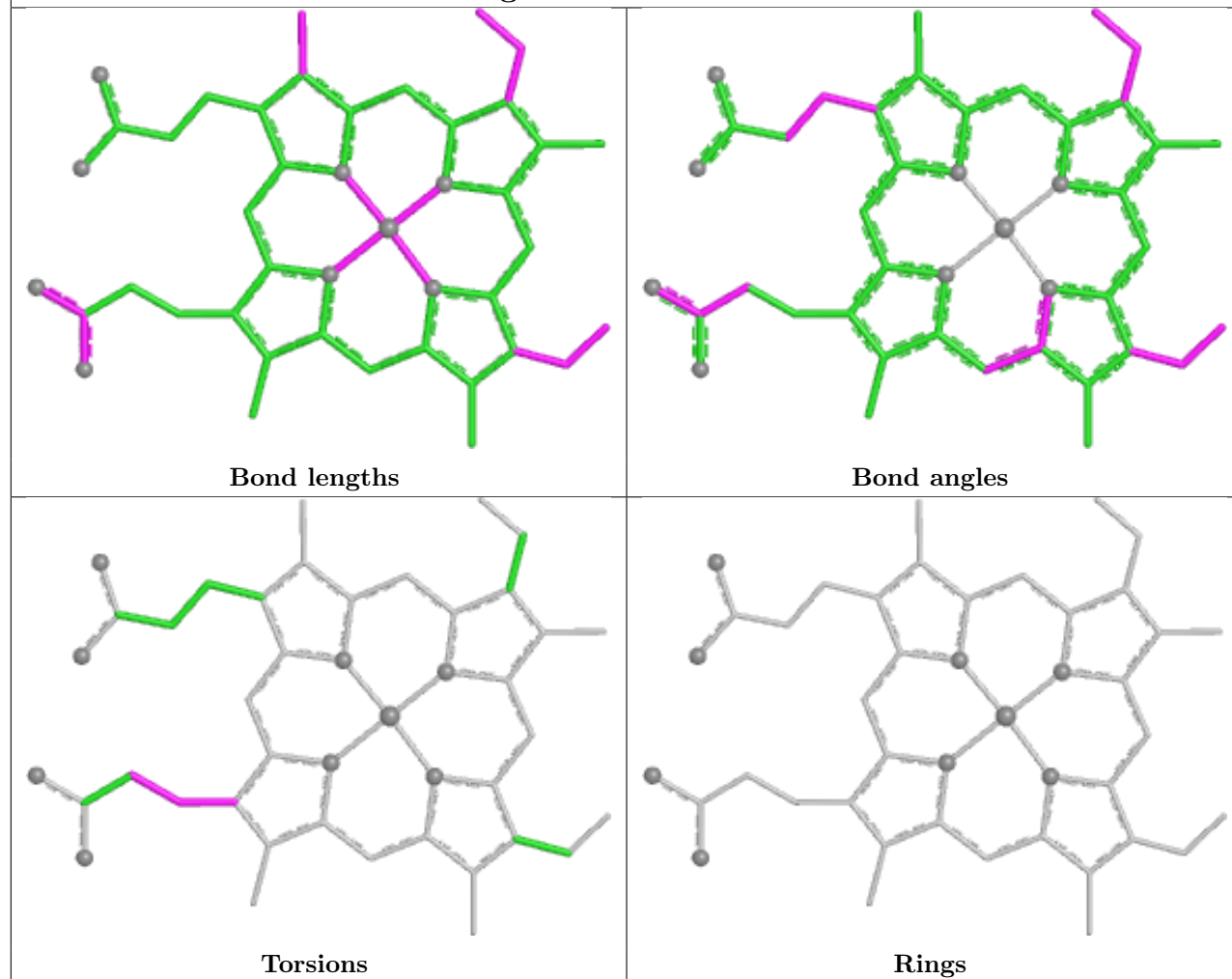


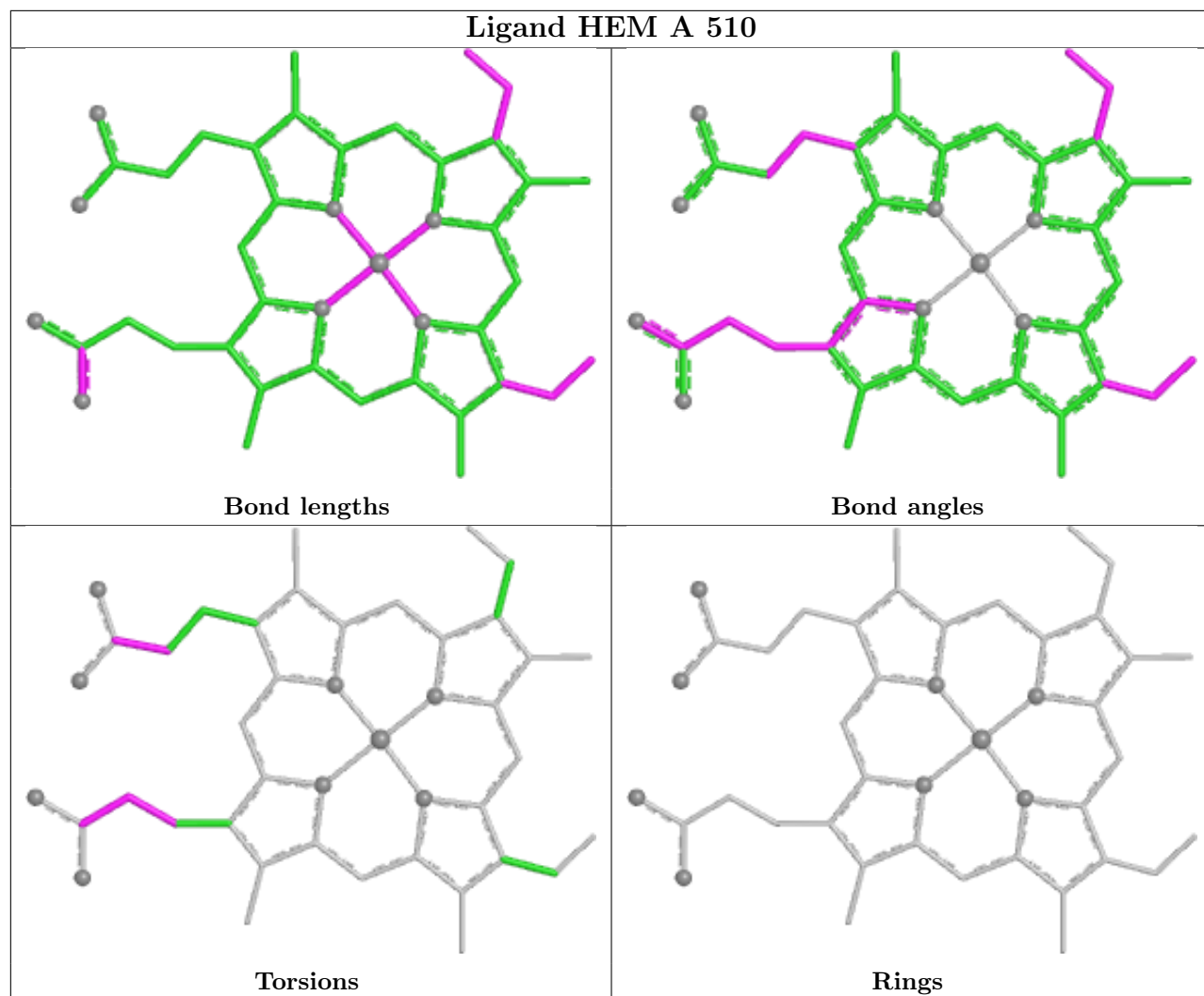
Torsions

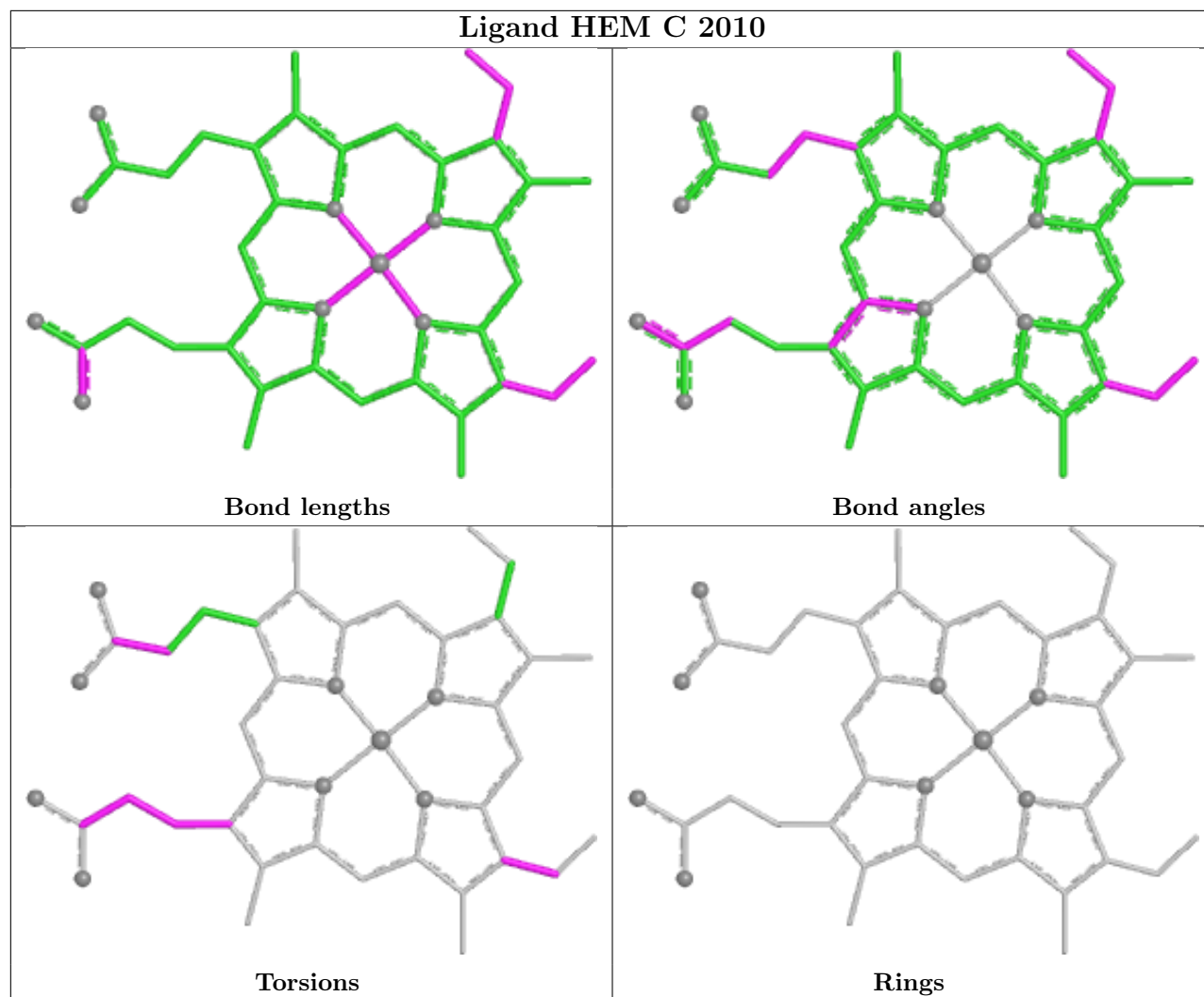


Rings

Ligand HEM B 1010







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.