



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 01:36 PM UTC

PDB ID : 1NSI / pdb_00001nsi
Title : HUMAN INDUCIBLE NITRIC OXIDE SYNTHASE, ZN-BOUND, L-ARG COMPLEX
Authors : Li, H.; Raman, C.S.; Glaser, C.B.; Blasko, E.; Young, T.A.; Parkinson, J.F.; Whitlow, M.; Poulos, T.L.
Deposited on : 1999-01-10
Resolution : 2.55 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

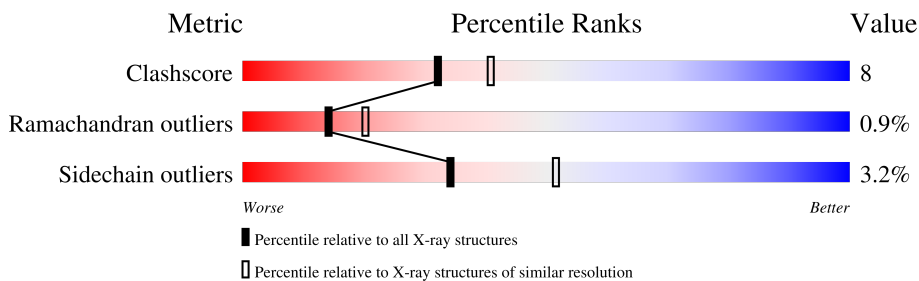
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.55 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	1120 (2.54-2.54)
Ramachandran outliers	187476	1106 (2.54-2.54)
Sidechain outliers	187428	1106 (2.54-2.54)

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	431	 79% 16% . .
1	B	431	 76% 18% . .
1	C	431	 78% 17% . .
1	D	431	 70% 25% . .

2 Entry composition [i](#)

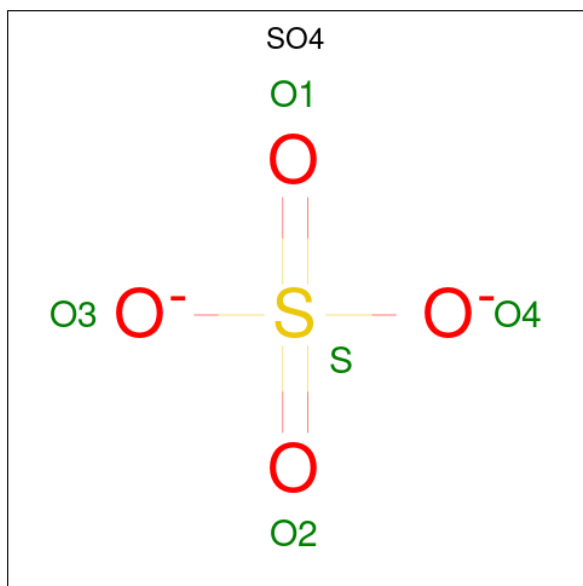
There are 8 unique types of molecules in this entry. The entry contains 14330 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 PROTEIN (NITRIC OXIDE SYNTHASE).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	420	Total	C	N	O	S	0	0	0
			3419	2186	599	612	22			
1	B	420	Total	C	N	O	S	0	0	0
			3419	2186	599	612	22			
1	C	420	Total	C	N	O	S	0	0	0
			3419	2186	599	612	22			
1	D	420	Total	C	N	O	S	0	0	0
			3419	2186	599	612	22			

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		

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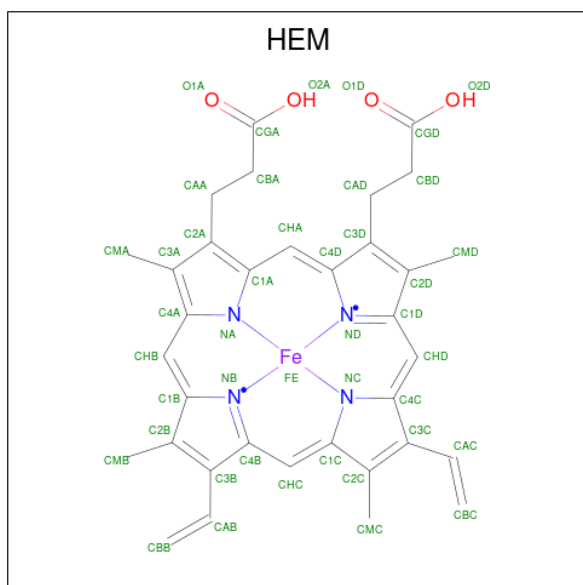
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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is ZINC ION (CCD ID: ZN) (formula: Zn).

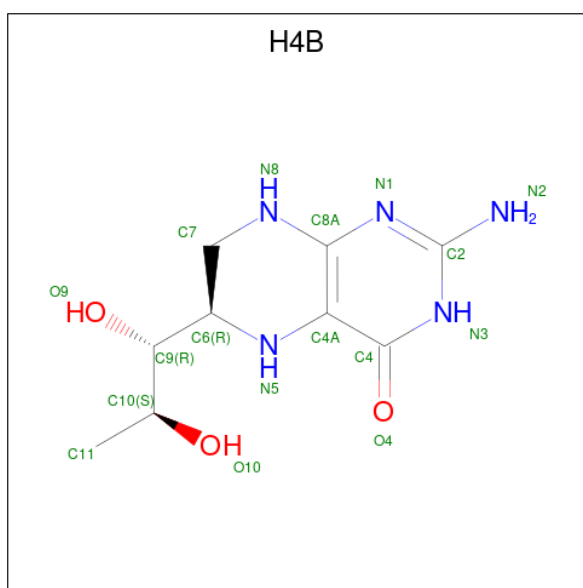
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Zn	0	0
			1	1		
3	C	1	Total	Zn	0	0
			1	1		

- Molecule 4 is PROTOPORPHYRIN IX CONTAINING FE (CCD ID: HEM) (formula: C₃₄H₃₂FeN₄O₄).



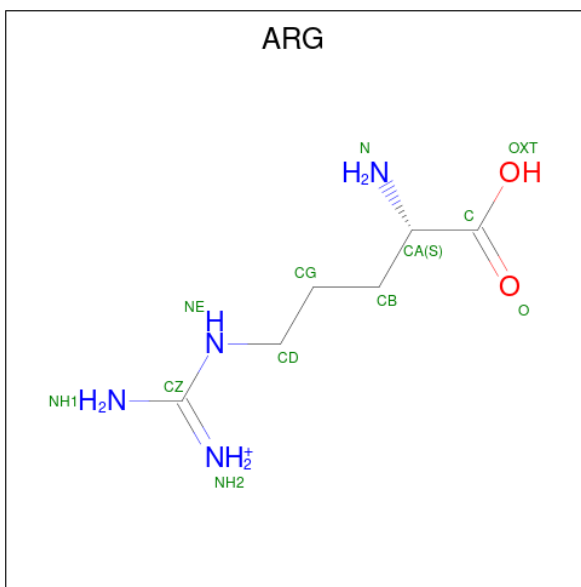
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	Fe	N	O	
			43	34	1	4	4	
4	B	1	Total	C	Fe	N	O	
			43	34	1	4	4	
4	C	1	Total	C	Fe	N	O	
			43	34	1	4	4	
4	D	1	Total	C	Fe	N	O	
			43	34	1	4	4	

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



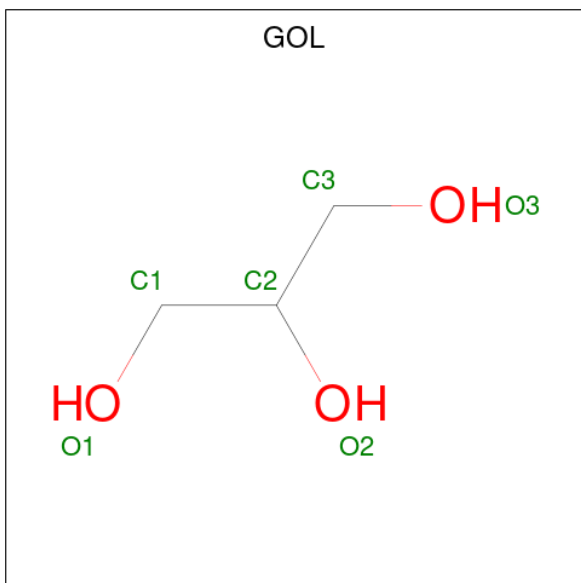
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	O		
			17	9	5	3	0	0
5	B	1	Total	C	N	O		
			17	9	5	3	0	0
5	C	1	Total	C	N	O		
			17	9	5	3	0	0
5	D	1	Total	C	N	O		
			17	9	5	3	0	0

- Molecule 6 is ARGinine (CCD ID: ARG) (formula: $C_6H_{15}N_4O_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	A	1	Total	C	N	O	0	0
			12	6	4	2		
6	B	1	Total	C	N	O	0	0
			12	6	4	2		
6	C	1	Total	C	N	O	0	0
			12	6	4	2		
6	D	1	Total	C	N	O	0	0
			12	6	4	2		

- Molecule 7 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



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

- Molecule 8 is water.

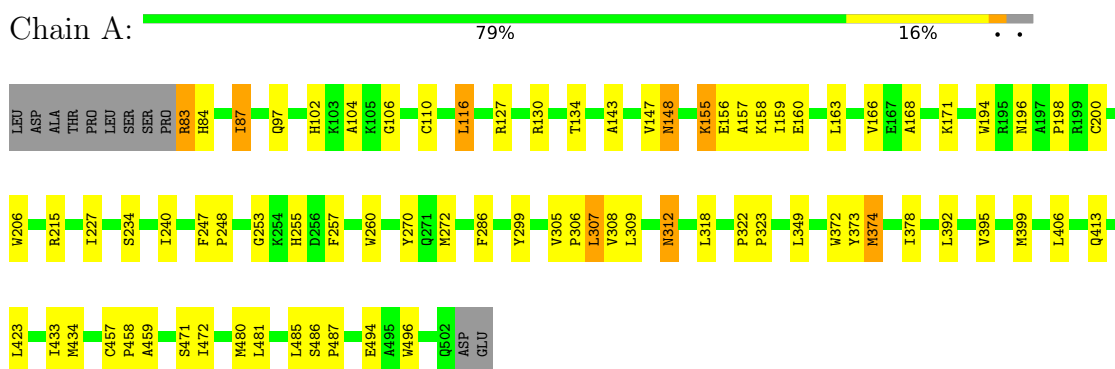
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	92	Total	O	0	0
			92	92		
8	B	77	Total	O	0	0
			77	77		
8	C	88	Total	O	0	0
			88	88		
8	D	38	Total	O	0	0
			38	38		

3 Residue-property plots

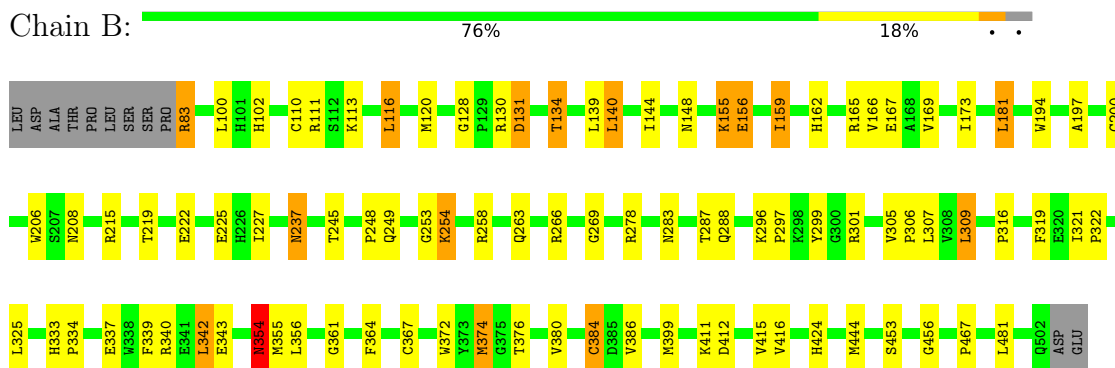
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded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

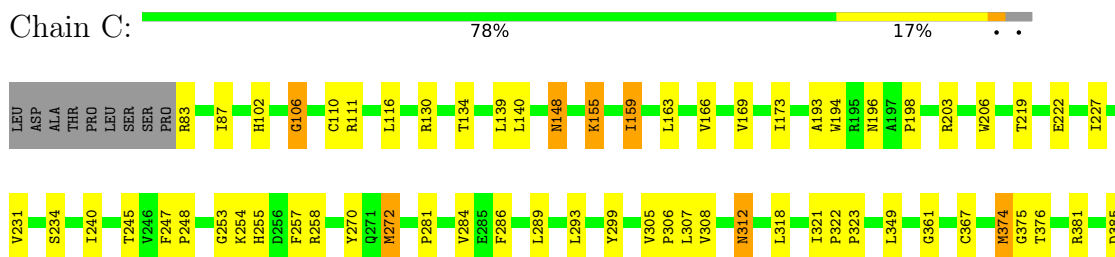
• Molecule 1: PROTEIN (NITRIC OXIDE SYNTHASE)



• Molecule 1: PROTEIN (NITRIC OXIDE SYNTHASE)

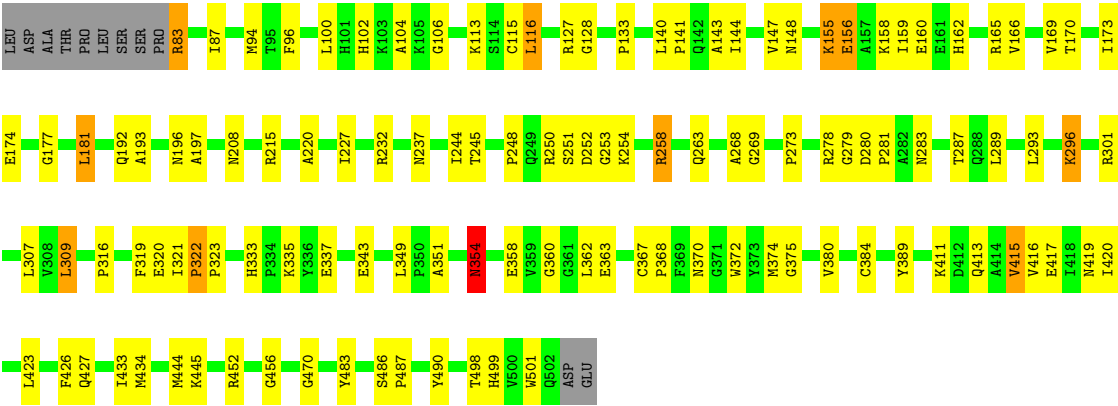


• Molecule 1: PROTEIN (NITRIC OXIDE SYNTHASE)





● Molecule 1: PROTEIN (NITRIC OXIDE SYNTHASE)



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	187.35Å 187.35Å 227.49Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.55	Depositor
% Data completeness (in resolution range)	80.5 (30.00-2.55)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	0.05	Depositor
Refinement program	X-PLOR 3.8	Depositor
R, R_{free}	0.209 , 0.243	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	14330	wwPDB-VP
Average B, all atoms (Å ²)	55.0	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, SO4, H4B, HEM, GOL

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	0.43	0/3518	0.93	8/4774 (0.2%)
1	B	0.44	0/3518	0.96	15/4774 (0.3%)
1	C	0.43	0/3518	0.92	10/4774 (0.2%)
1	D	0.45	0/3518	0.99	20/4774 (0.4%)
All	All	0.44	0/14072	0.95	53/19096 (0.3%)

There are no bond length outliers.

All (53) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	253	GLY	N-CA-C	-8.54	103.23	115.27
1	C	253	GLY	N-CA-C	-8.44	103.37	115.27
1	A	253	GLY	N-CA-C	-7.68	103.89	114.64
1	D	253	GLY	N-CA-C	-7.26	104.55	115.00
1	D	501	TRP	N-CA-C	7.09	119.37	109.15
1	A	200	CYS	CA-CB-SG	-7.04	98.22	114.40
1	B	384	CYS	N-CA-C	6.62	121.44	113.23
1	B	258	ARG	N-CA-C	6.53	119.24	108.99
1	D	258	ARG	N-CA-C	6.48	119.57	109.52
1	D	349	LEU	N-CA-C	6.41	120.23	108.69
1	C	374	MET	N-CA-C	-6.36	98.54	108.90
1	B	322	PRO	CA-C-N	5.96	126.39	119.47
1	B	322	PRO	C-N-CA	5.96	126.39	119.47
1	A	309	LEU	N-CA-C	5.89	119.42	109.76
1	C	272	MET	CA-C-N	5.89	125.76	119.28
1	C	272	MET	C-N-CA	5.89	125.76	119.28
1	C	254	LYS	N-CA-C	-5.88	106.44	113.97
1	A	392	LEU	N-CA-C	5.84	117.32	111.07
1	B	206	TRP	N-CA-C	5.78	118.04	111.11
1	C	203	ARG	N-CA-C	5.74	119.76	112.87

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	355	MET	N-CA-C	5.72	118.88	110.59
1	A	373	TYR	N-CA-C	5.67	118.01	110.53
1	D	322	PRO	CA-C-N	5.64	126.01	119.47
1	D	322	PRO	C-N-CA	5.64	126.01	119.47
1	B	374	MET	N-CA-C	-5.59	100.07	109.07
1	B	197	ALA	CA-C-N	5.52	125.23	119.82
1	B	197	ALA	C-N-CA	5.52	125.23	119.82
1	A	378	ILE	CB-CA-C	-5.50	104.84	112.04
1	A	349	LEU	N-CA-C	5.48	118.55	108.69
1	C	408	SER	N-CA-C	-5.47	106.61	113.28
1	D	128	GLY	CA-C-N	5.47	125.67	120.31
1	D	128	GLY	C-N-CA	5.47	125.67	120.31
1	B	194	TRP	N-CA-C	-5.41	105.30	111.14
1	C	479	GLU	N-CA-C	-5.39	101.69	110.20
1	D	197	ALA	CA-C-N	5.39	125.06	119.56
1	D	197	ALA	C-N-CA	5.39	125.06	119.56
1	D	384	CYS	N-CA-C	5.37	119.98	113.38
1	B	120	MET	N-CA-C	5.30	117.48	111.02
1	D	354	ASN	N-CA-C	5.29	119.60	113.20
1	D	389	TYR	N-CA-C	5.29	117.79	111.71
1	D	420	ILE	N-CA-C	-5.28	105.38	110.72
1	B	254	LYS	N-CA-C	-5.19	106.34	113.30
1	D	415	VAL	CB-CA-C	-5.18	105.24	111.88
1	B	354	ASN	N-CA-C	5.18	120.52	113.37
1	B	131	ASP	N-CA-C	-5.17	106.08	114.09
1	C	349	LEU	N-CA-C	5.15	117.97	108.69
1	A	374	MET	N-CA-C	-5.13	100.16	108.52
1	D	220	ALA	N-CA-C	-5.12	105.78	111.36
1	D	280	ASP	CA-C-N	5.08	124.74	119.56
1	D	280	ASP	C-N-CA	5.08	124.74	119.56
1	C	258	ARG	N-CA-C	5.06	116.75	108.76
1	D	470	GLY	N-CA-C	5.00	118.31	112.50
1	D	254	LYS	N-CA-C	-5.00	107.22	113.72

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	3419	0	3326	50	0
1	B	3419	0	3326	63	0
1	C	3419	0	3326	54	0
1	D	3419	0	3326	67	0
2	A	15	0	0	0	0
2	B	10	0	0	0	0
2	C	10	0	0	0	0
2	D	10	0	0	1	0
3	A	1	0	0	0	0
3	C	1	0	0	0	0
4	A	43	0	30	1	0
4	B	43	0	30	2	0
4	C	43	0	30	0	0
4	D	43	0	30	3	0
5	A	17	0	15	0	0
5	B	17	0	15	0	0
5	C	17	0	15	1	0
5	D	17	0	15	0	0
6	A	12	0	12	0	0
6	B	12	0	12	0	0
6	C	12	0	12	0	0
6	D	12	0	12	0	0
7	A	6	0	8	0	0
7	B	6	0	8	0	0
7	C	6	0	8	0	0
7	D	6	0	8	3	0
8	A	92	0	0	1	0
8	B	77	0	0	0	0
8	C	88	0	0	0	0
8	D	38	0	0	0	0
All	All	14330	0	13564	222	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 8.

All (222) 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:481:LEU:HD22	1:B:116:LEU:HD22	1.54	0.90

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:270:TYR:HB2	1:C:272:MET:HE3	1.57	0.86
1:C:481:LEU:HD22	1:D:116:LEU:HD22	1.56	0.86
1:A:83:ARG:HH11	1:A:83:ARG:HB2	1.42	0.84
1:D:411:LYS:O	1:D:415:VAL:HG23	1.81	0.80
1:A:116:LEU:HD22	1:B:481:LEU:HG	1.68	0.76
1:D:83:ARG:HH11	1:D:83:ARG:HB2	1.49	0.75
1:C:227:ILE:HG21	1:C:307:LEU:HD21	1.74	0.69
1:C:257:PHE:CE1	1:C:312:ASN:HB3	2.27	0.68
1:B:269:GLY:H	1:B:287:THR:HG21	1.59	0.68
1:A:272:MET:HA	1:A:272:MET:HE2	1.76	0.67
1:A:372:TRP:H	4:A:550:HEM:HAB	1.58	0.66
1:B:305:VAL:HG22	1:B:306:PRO:HD2	1.77	0.66
1:D:227:ILE:HG21	1:D:307:LEU:HD21	1.77	0.66
1:D:100:LEU:HB3	1:D:456:GLY:HA3	1.78	0.66
1:B:227:ILE:HG21	1:B:307:LEU:HD21	1.76	0.65
1:D:258:ARG:HH21	7:D:883:GOL:H2	1.60	0.65
1:B:411:LYS:O	1:B:415:VAL:HG23	1.97	0.64
1:D:140:LEU:O	1:D:144:ILE:HG12	1.98	0.64
1:A:83:ARG:HB2	1:A:83:ARG:NH1	2.13	0.64
1:B:144:ILE:HG22	1:B:148:ASN:HD21	1.63	0.63
1:D:143:ALA:O	1:D:147:VAL:HG23	1.99	0.63
1:D:354:ASN:HD22	1:D:354:ASN:H	1.46	0.62
1:A:84:HIS:HB3	1:A:97:GLN:NE2	2.15	0.62
1:C:255:HIS:HB3	1:C:312:ASN:HB2	1.82	0.62
1:B:283:ASN:O	1:B:287:THR:HG23	2.00	0.61
1:C:130:ARG:NH1	1:C:134:THR:HG22	2.17	0.60
1:B:354:ASN:N	1:B:354:ASN:HD22	2.00	0.60
1:A:83:ARG:HB3	1:A:102:HIS:NE2	2.18	0.59
1:D:170:THR:O	1:D:174:GLU:HB2	2.02	0.59
1:B:83:ARG:HB2	1:B:83:ARG:HH11	1.67	0.59
1:D:301:ARG:HH22	1:D:343:GLU:HB2	1.67	0.59
1:D:269:GLY:H	1:D:287:THR:HG21	1.67	0.59
1:B:111:ARG:HD2	1:B:111:ARG:N	2.17	0.59
1:A:257:PHE:CE1	1:A:312:ASN:HB3	2.38	0.58
1:A:227:ILE:HG21	1:A:307:LEU:HD21	1.84	0.58
1:B:215:ARG:O	1:B:248:PRO:HG3	2.03	0.57
1:C:412:ASP:O	1:C:416:VAL:HG23	2.03	0.57
1:B:140:LEU:HD22	1:B:173:ILE:HD12	1.87	0.56
1:A:168:ALA:O	1:A:171:LYS:HE2	2.05	0.56
1:B:83:ARG:HB3	1:B:102:HIS:NE2	2.20	0.56
1:C:305:VAL:HG13	1:C:306:PRO:HD2	1.86	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:309:LEU:HD23	1:D:319:PHE:CD2	2.40	0.56
1:B:372:TRP:H	4:B:550:HEM:HAB	1.71	0.56
1:D:162:HIS:O	1:D:166:VAL:HG23	2.07	0.55
1:A:472:ILE:HG13	1:B:415:VAL:HG21	1.88	0.55
1:B:130:ARG:CZ	1:B:134:THR:HG23	2.36	0.55
1:D:333:HIS:CE1	1:D:335:LYS:HG3	2.42	0.55
1:B:139:LEU:CD1	1:B:361:GLY:HA3	2.37	0.55
1:D:372:TRP:H	4:D:550:HEM:HAB	1.72	0.55
1:D:283:ASN:O	1:D:287:THR:HG23	2.07	0.54
1:B:100:LEU:HB3	1:B:456:GLY:HA3	1.88	0.54
1:B:181:LEU:CD2	1:B:181:LEU:H	2.20	0.54
1:A:83:ARG:HB3	1:A:102:HIS:CE1	2.42	0.54
1:D:232:ARG:HH12	1:D:237:ASN:HD21	1.56	0.54
1:D:144:ILE:HG22	1:D:148:ASN:HD21	1.73	0.54
1:A:196:ASN:O	1:A:198:PRO:HD3	2.08	0.54
1:C:83:ARG:HB3	1:C:102:HIS:NE2	2.23	0.53
1:A:308:VAL:HG22	1:A:318:LEU:HD22	1.90	0.53
1:A:471:SER:OG	1:B:467:PRO:HB2	2.09	0.53
1:C:374:MET:HE3	1:C:376:THR:OG1	2.10	0.52
1:D:354:ASN:HD22	1:D:354:ASN:N	2.08	0.52
1:D:445:LYS:HZ1	1:D:452:ARG:HH22	1.57	0.52
1:D:309:LEU:O	1:D:316:PRO:HA	2.09	0.52
1:C:270:TYR:CB	1:C:272:MET:HE3	2.35	0.52
1:D:94:MET:HE3	1:D:96:PHE:CZ	2.45	0.52
1:D:380:VAL:HG22	1:D:419:ASN:HD21	1.74	0.52
1:A:413:GLN:HG2	1:B:416:VAL:HG11	1.92	0.51
1:C:194:TRP:CE3	1:C:206:TRP:HA	2.45	0.51
1:C:169:VAL:O	1:C:173:ILE:HG13	2.11	0.51
1:D:370:ASN:O	4:D:550:HEM:HMC3	2.11	0.51
1:A:286:PHE:CE2	1:A:305:VAL:HG11	2.46	0.50
1:D:367:CYS:N	1:D:368:PRO:HD3	2.25	0.50
1:A:286:PHE:HE2	1:A:305:VAL:HG11	1.77	0.50
8:A:1020:HOH:O	1:B:386:VAL:HG22	2.11	0.50
1:D:165:ARG:O	1:D:169:VAL:HG23	2.11	0.50
1:A:143:ALA:O	1:A:147:VAL:HG23	2.12	0.50
1:A:194:TRP:CE3	1:A:206:TRP:HA	2.47	0.50
1:B:165:ARG:O	1:B:169:VAL:HG23	2.12	0.49
1:B:131:ASP:O	1:B:254:LYS:HE3	2.12	0.49
1:B:339:PHE:O	1:B:342:LEU:HB2	2.11	0.49
1:B:140:LEU:HD12	1:C:140:LEU:CD1	2.43	0.49
1:C:130:ARG:CZ	1:C:134:THR:HG22	2.43	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:305:VAL:CG2	1:B:306:PRO:HD2	2.41	0.49
1:C:413:GLN:HG2	1:D:416:VAL:HG11	1.95	0.49
1:C:245:THR:O	1:C:367:CYS:HA	2.13	0.49
1:D:83:ARG:HB3	1:D:102:HIS:NE2	2.28	0.49
1:C:381:ARG:HH12	5:C:602:H4B:C4	2.26	0.48
1:A:106:GLY:HA2	1:B:113:LYS:HA	1.95	0.48
1:C:486:SER:HA	1:C:487:PRO:C	2.39	0.48
1:A:155:LYS:CD	1:A:155:LYS:H	2.27	0.48
1:D:358:GLU:O	1:D:487:PRO:HA	2.13	0.48
1:D:155:LYS:HE3	1:D:155:LYS:H	1.77	0.48
1:B:374:MET:HE3	1:B:376:THR:OG1	2.13	0.48
1:C:248:PRO:HG2	1:C:257:PHE:CD1	2.49	0.48
1:D:140:LEU:HD11	1:D:170:THR:HG23	1.94	0.48
1:D:423:LEU:O	1:D:427:GLN:HG3	2.14	0.47
1:A:163:LEU:HA	1:A:166:VAL:HG12	1.96	0.47
1:B:412:ASP:O	1:B:416:VAL:HG23	2.14	0.47
1:C:155:LYS:H	1:C:155:LYS:HD3	1.80	0.47
1:B:333:HIS:ND1	1:B:334:PRO:HD2	2.29	0.47
1:A:110:CYS:HB3	1:B:110:CYS:HB3	1.96	0.47
1:D:289:LEU:O	1:D:293:LEU:HG	2.14	0.47
1:A:255:HIS:HB3	1:A:312:ASN:HB2	1.96	0.47
1:D:445:LYS:NZ	1:D:452:ARG:NH2	2.62	0.47
1:C:293:LEU:HD12	1:C:318:LEU:HD21	1.97	0.47
1:C:385:ASP:HB2	1:C:388:ARG:HG2	1.96	0.47
1:A:234:SER:O	1:A:240:ILE:HA	2.15	0.47
1:D:140:LEU:N	1:D:141:PRO:CD	2.78	0.47
1:B:110:CYS:C	1:B:111:ARG:HD2	2.40	0.47
1:A:305:VAL:HG13	1:A:306:PRO:HD2	1.97	0.46
1:B:222:GLU:HA	1:B:225:GLU:OE1	2.15	0.46
1:A:158:LYS:O	1:A:160:GLU:N	2.48	0.46
1:D:158:LYS:O	1:D:160:GLU:N	2.48	0.46
1:D:173:ILE:O	1:D:177:GLY:HA2	2.15	0.46
1:C:196:ASN:O	1:C:198:PRO:HD3	2.16	0.46
1:D:374:MET:HA	1:D:434:MET:O	2.15	0.46
1:C:374:MET:HA	1:C:434:MET:O	2.16	0.46
1:B:380:VAL:O	1:B:384:CYS:HB2	2.16	0.45
1:C:272:MET:HE1	1:C:299:TYR:CD1	2.51	0.45
1:A:215:ARG:HA	1:A:248:PRO:HD3	1.99	0.45
1:A:247:PHE:HB3	1:A:248:PRO:HD2	1.98	0.45
1:A:395:VAL:O	1:A:399:MET:HG3	2.16	0.45
1:B:263:GLN:HB2	1:B:266:ARG:HD2	1.97	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:308:VAL:HG22	1:A:318:LEU:CD2	2.46	0.45
1:B:340:ARG:HE	1:B:340:ARG:HB3	1.55	0.45
1:C:286:PHE:CE2	1:C:305:VAL:HG11	2.52	0.45
1:D:193:ALA:HB2	1:D:487:PRO:HB2	1.99	0.45
1:B:237:ASN:N	1:B:237:ASN:HD22	2.14	0.45
1:C:281:PRO:HA	1:C:284:VAL:HG23	1.99	0.45
1:C:385:ASP:HB2	1:C:388:ARG:CG	2.47	0.45
1:D:498:THR:OG1	7:D:883:GOL:H32	2.17	0.45
1:C:247:PHE:HB3	1:C:248:PRO:HD2	1.99	0.44
1:B:301:ARG:HH22	1:B:343:GLU:HB2	1.83	0.44
1:B:356:LEU:C	1:B:356:LEU:HD23	2.42	0.44
1:D:87:ILE:HG13	1:D:96:PHE:HB2	1.98	0.44
1:A:104:ALA:HA	1:A:485:LEU:HD23	2.00	0.44
1:A:260:TRP:CZ3	1:A:496:TRP:HB3	2.52	0.44
1:A:148:ASN:N	1:A:148:ASN:HD22	2.16	0.44
1:B:334:PRO:HG3	1:B:424:HIS:CG	2.52	0.44
1:A:486:SER:HA	1:A:487:PRO:C	2.43	0.44
1:C:227:ILE:O	1:C:231:VAL:HG23	2.17	0.44
1:D:301:ARG:NH2	1:D:343:GLU:HB2	2.33	0.44
1:D:322:PRO:HA	1:D:323:PRO:HD3	1.88	0.44
1:D:263:GLN:HA	1:D:351:ALA:O	2.18	0.44
1:C:155:LYS:H	1:C:155:LYS:CD	2.31	0.43
1:A:272:MET:HE1	1:A:299:TYR:CD1	2.53	0.43
1:C:308:VAL:HG22	1:C:318:LEU:HD22	2.01	0.43
1:C:106:GLY:HA2	1:D:113:LYS:HA	1.99	0.43
1:B:245:THR:O	1:B:367:CYS:HA	2.19	0.43
1:C:451:TYR:CE1	1:C:457:CYS:HB3	2.53	0.43
1:A:270:TYR:CB	1:A:272:MET:HE3	2.48	0.43
1:B:296:LYS:HA	1:B:297:PRO:HD2	1.89	0.43
1:D:250:ARG:HD2	1:D:363:GLU:OE1	2.18	0.43
1:A:130:ARG:NH1	1:A:134:THR:HG22	2.34	0.43
1:A:155:LYS:H	1:A:155:LYS:HE3	1.83	0.43
1:D:307:LEU:HD13	1:D:321:ILE:HD11	2.00	0.43
1:B:307:LEU:HD22	1:B:321:ILE:HG12	2.01	0.43
1:B:354:ASN:HD22	1:B:354:ASN:H	1.67	0.43
1:D:499:HIS:HA	7:D:883:GOL:O3	2.19	0.43
1:B:140:LEU:O	1:B:144:ILE:HG12	2.18	0.43
1:D:87:ILE:CG1	1:D:96:PHE:HB2	2.49	0.43
1:A:156:GLU:HB3	1:A:157:ALA:H	1.66	0.42
1:A:127:ARG:HD3	1:A:127:ARG:HA	1.85	0.42
1:A:248:PRO:HG2	1:A:257:PHE:CG	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:459:ALA:HB3	1:A:480:MET:HB2	2.00	0.42
1:B:181:LEU:CD2	1:B:181:LEU:N	2.83	0.42
1:B:342:LEU:HD12	1:B:342:LEU:HA	1.87	0.42
1:D:296:LYS:HD3	1:D:296:LYS:N	2.35	0.42
1:A:270:TYR:HB3	1:A:272:MET:HE3	2.01	0.42
1:B:200:CYS:HB2	4:B:550:HEM:ND	2.35	0.42
1:C:148:ASN:N	1:C:148:ASN:HD22	2.16	0.42
1:C:375:GLY:HA3	1:C:423:LEU:HD11	2.02	0.42
1:D:244:ILE:HG12	1:D:245:THR:N	2.33	0.42
1:D:426:PHE:CD2	1:D:433:ILE:HB	2.55	0.42
1:A:87:ILE:HD13	1:A:87:ILE:HA	1.77	0.42
1:B:159:ILE:HG22	1:C:159:ILE:HG22	2.01	0.42
1:B:249:GLN:HB3	1:B:364:PHE:CE2	2.54	0.42
1:C:193:ALA:HB2	1:C:487:PRO:HB2	2.01	0.42
1:D:268:ALA:HB1	1:D:287:THR:HG21	2.01	0.42
1:B:155:LYS:H	1:B:155:LYS:HE3	1.85	0.42
1:B:181:LEU:N	1:B:181:LEU:HD23	2.35	0.42
1:C:270:TYR:CE1	1:C:299:TYR:HA	2.55	0.42
1:D:133:PRO:HB3	1:D:251:SER:O	2.19	0.42
1:C:289:LEU:O	1:C:293:LEU:HG	2.20	0.42
1:D:155:LYS:NZ	1:D:156:GLU:HG3	2.35	0.42
1:D:337:GLU:HB3	2:D:923:SO4:O2	2.19	0.42
1:B:155:LYS:NZ	1:B:156:GLU:HG3	2.35	0.41
1:B:155:LYS:HZ1	1:B:156:GLU:HG3	1.85	0.41
1:C:412:ASP:O	1:C:415:VAL:HG12	2.20	0.41
1:B:309:LEU:HD23	1:B:319:PHE:CD1	2.55	0.41
1:D:215:ARG:O	1:D:248:PRO:HG3	2.20	0.41
1:B:162:HIS:O	1:B:166:VAL:HG23	2.20	0.41
1:C:234:SER:O	1:C:240:ILE:HA	2.20	0.41
1:C:110:CYS:O	1:C:111:ARG:HD3	2.20	0.41
1:A:374:MET:HA	1:A:434:MET:O	2.21	0.41
1:C:322:PRO:HA	1:C:323:PRO:HD2	1.91	0.41
1:D:181:LEU:HD22	1:D:362:LEU:CD1	2.51	0.41
1:C:367:CYS:O	1:C:367:CYS:SG	2.79	0.41
1:C:437:HIS:O	1:C:441:GLU:HG3	2.21	0.41
1:D:115:CYS:C	1:D:116:LEU:HD23	2.46	0.41
1:A:457:CYS:HA	1:A:458:PRO:HD2	1.96	0.41
1:C:219:THR:OG1	1:C:222:GLU:HG3	2.21	0.41
4:D:550:HEM:HBC2	4:D:550:HEM:HMC2	2.02	0.41
1:D:127:ARG:HB2	1:D:490:TYR:CB	2.50	0.41
1:B:144:ILE:HG22	1:B:148:ASN:ND2	2.32	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:309:LEU:O	1:B:316:PRO:HA	2.21	0.41
1:B:354:ASN:N	1:B:354:ASN:ND2	2.65	0.41
1:C:307:LEU:HD13	1:C:321:ILE:HD11	2.02	0.41
1:D:486:SER:HA	1:D:487:PRO:C	2.45	0.41
1:B:354:ASN:H	1:B:354:ASN:ND2	2.19	0.41
1:C:472:ILE:HG13	1:D:415:VAL:HG21	2.03	0.41
1:D:192:GLN:HG3	1:D:196:ASN:HD21	1.86	0.41
1:A:322:PRO:HA	1:A:323:PRO:HD2	1.86	0.40
1:D:279:GLY:O	1:D:281:PRO:HD3	2.21	0.40
1:C:83:ARG:HG2	1:C:83:ARG:NH1	2.36	0.40
1:C:139:LEU:HD12	1:C:361:GLY:HA3	2.02	0.40
1:C:286:PHE:HE2	1:C:305:VAL:HG11	1.87	0.40
1:D:413:GLN:O	1:D:417:GLU:HG2	2.21	0.40
1:B:399:MET:HE2	1:B:399:MET:HB3	1.96	0.40
1:D:104:ALA:CB	1:D:483:TYR:HB2	2.51	0.40
1:C:163:LEU:HA	1:C:166:VAL:HG12	2.04	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	418/431 (97%)	388 (93%)	28 (7%)	2 (0%)	24	34
1	B	418/431 (97%)	383 (92%)	30 (7%)	5 (1%)	10	13
1	C	418/431 (97%)	386 (92%)	29 (7%)	3 (1%)	18	25
1	D	418/431 (97%)	381 (91%)	32 (8%)	5 (1%)	10	13
All	All	1672/1724 (97%)	1538 (92%)	119 (7%)	15 (1%)	14	20

All (15) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	159	ILE
1	B	299	TYR
1	D	159	ILE
1	B	159	ILE
1	C	159	ILE
1	C	312	ASN
1	B	156	GLU
1	B	453	SER
1	A	312	ASN
1	B	128	GLY
1	D	156	GLU
1	C	106	GLY
1	D	375	GLY
1	D	106	GLY
1	D	360	GLY

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	366/376 (97%)	356 (97%)	10 (3%)	39	58
1	B	366/376 (97%)	348 (95%)	18 (5%)	22	34
1	C	366/376 (97%)	360 (98%)	6 (2%)	55	73
1	D	366/376 (97%)	353 (96%)	13 (4%)	31	46
All	All	1464/1504 (97%)	1417 (97%)	47 (3%)	34	51

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

Mol	Chain	Res	Type
1	A	83	ARG
1	A	87	ILE
1	A	116	LEU
1	A	148	ASN
1	A	155	LYS
1	A	307	LEU
1	A	406	LEU

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Mol	Chain	Res	Type
1	A	423	LEU
1	A	433	ILE
1	A	494	GLU
1	B	83	ARG
1	B	116	LEU
1	B	134	THR
1	B	140	LEU
1	B	155	LYS
1	B	167	GLU
1	B	181	LEU
1	B	208	ASN
1	B	219	THR
1	B	237	ASN
1	B	278	ARG
1	B	288	GLN
1	B	309	LEU
1	B	325	LEU
1	B	337	GLU
1	B	342	LEU
1	B	354	ASN
1	B	444	MET
1	C	87	ILE
1	C	116	LEU
1	C	148	ASN
1	C	155	LYS
1	C	433	ILE
1	C	494	GLU
1	D	83	ARG
1	D	116	LEU
1	D	155	LYS
1	D	181	LEU
1	D	208	ASN
1	D	252	ASP
1	D	273	PRO
1	D	278	ARG
1	D	296	LYS
1	D	309	LEU
1	D	320	GLU
1	D	354	ASN
1	D	444	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (44) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	97	GLN
1	A	148	ASN
1	A	162	HIS
1	A	205	GLN
1	A	390	ASN
1	A	424	HIS
1	A	427	GLN
1	A	482	ASN
1	B	148	ASN
1	B	192	GLN
1	B	205	GLN
1	B	237	ASN
1	B	263	GLN
1	B	271	GLN
1	B	288	GLN
1	B	312	ASN
1	B	354	ASN
1	B	390	ASN
1	B	424	HIS
1	B	427	GLN
1	B	502	GLN
1	C	97	GLN
1	C	148	ASN
1	C	162	HIS
1	C	192	GLN
1	C	205	GLN
1	C	210	GLN
1	C	390	ASN
1	C	413	GLN
1	C	427	GLN
1	C	482	ASN
1	D	148	ASN
1	D	192	GLN
1	D	205	GLN
1	D	208	ASN
1	D	237	ASN
1	D	249	GLN
1	D	263	GLN
1	D	271	GLN
1	D	312	ASN
1	D	354	ASN
1	D	370	ASN
1	D	390	ASN

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Mol	Chain	Res	Type
1	D	427	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](#)

Of 27 ligands modelled in this entry, 2 are monoatomic - leaving 25 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	$\# Z > 2$
2	SO4	A	910	-	4,4,4	0.62	0	6,6,6	0.34	0
6	ARG	A	700	-	10,11,11	0.65	0	9,13,13	0.38	0
6	ARG	D	703	-	10,11,11	0.65	0	9,13,13	0.35	0
2	SO4	A	930	-	4,4,4	0.67	0	6,6,6	0.32	0
4	HEM	A	550	1	50,50,50	1.12	2 (4%)	67,82,82	1.16	4 (5%)
2	SO4	C	912	-	4,4,4	0.63	0	6,6,6	0.22	0
6	ARG	B	701	-	10,11,11	0.71	0	9,13,13	0.31	0
5	H4B	A	600	-	17,18,18	1.14	2 (11%)	14,26,26	1.85	3 (21%)
5	H4B	B	601	-	17,18,18	1.21	2 (11%)	14,26,26	1.87	3 (21%)
7	GOL	A	880	-	5,5,5	0.40	0	5,5,5	0.31	0
2	SO4	D	913	-	4,4,4	0.71	0	6,6,6	0.23	0
4	HEM	B	550	1	50,50,50	1.08	3 (6%)	67,82,82	1.34	8 (11%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
7	GOL	C	882	-	5,5,5	0.43	0	5,5,5	0.31	0
4	HEM	C	550	1	50,50,50	1.31	6 (12%)	67,82,82	1.35	8 (11%)
7	GOL	B	881	-	5,5,5	0.33	0	5,5,5	0.30	0
6	ARG	C	702	-	10,11,11	0.70	0	9,13,13	0.40	0
7	GOL	D	883	-	5,5,5	0.40	0	5,5,5	0.24	0
2	SO4	D	923	-	4,4,4	0.72	0	6,6,6	0.33	0
2	SO4	A	920	-	4,4,4	0.73	0	6,6,6	0.30	0
2	SO4	C	922	-	4,4,4	0.69	0	6,6,6	0.29	0
5	H4B	C	602	-	17,18,18	1.18	2 (11%)	14,26,26	1.89	3 (21%)
4	HEM	D	550	1	50,50,50	1.16	3 (6%)	67,82,82	1.33	5 (7%)
5	H4B	D	603	-	17,18,18	1.27	2 (11%)	14,26,26	1.84	3 (21%)
2	SO4	B	911	-	4,4,4	0.70	0	6,6,6	0.36	0
2	SO4	B	921	-	4,4,4	0.66	0	6,6,6	0.46	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	H4B	A	600	-	-	0/8/17/17	0/2/2/2
5	H4B	B	601	-	-	0/8/17/17	0/2/2/2
6	ARG	A	700	-	-	0/11/11/11	-
6	ARG	D	703	-	-	0/11/11/11	-
7	GOL	A	880	-	-	4/4/4/4	-
4	HEM	B	550	1	-	4/14/54/54	-
7	GOL	C	882	-	-	1/4/4/4	-
5	H4B	C	602	-	-	0/8/17/17	0/2/2/2
4	HEM	C	550	1	-	2/14/54/54	-
7	GOL	B	881	-	-	4/4/4/4	-
4	HEM	A	550	1	-	2/14/54/54	-
4	HEM	D	550	1	-	1/14/54/54	-
6	ARG	B	701	-	-	0/11/11/11	-
5	H4B	D	603	-	-	0/8/17/17	0/2/2/2
6	ARG	C	702	-	-	0/11/11/11	-
7	GOL	D	883	-	-	4/4/4/4	-

All (22) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	550	HEM	CAB-C3B	-3.85	1.37	1.47
5	D	603	H4B	C4A-N5	3.42	1.45	1.37
4	A	550	HEM	CAC-C3C	-3.24	1.38	1.47
5	C	602	H4B	C4A-N5	3.18	1.45	1.37
5	B	601	H4B	C4A-N5	3.16	1.45	1.37
4	D	550	HEM	CAB-C3B	-3.14	1.39	1.47
5	A	600	H4B	C4A-N5	3.14	1.45	1.37
4	C	550	HEM	CAC-C3C	-3.02	1.39	1.47
4	C	550	HEM	C4D-ND	-2.98	1.35	1.40
4	B	550	HEM	CAB-C3B	-2.96	1.39	1.47
4	C	550	HEM	CAB-C3B	-2.96	1.39	1.47
4	C	550	HEM	C3D-C2D	-2.91	1.30	1.36
5	A	600	H4B	C6-N5	2.50	1.51	1.47
5	C	602	H4B	C6-N5	2.47	1.51	1.47
4	D	550	HEM	CAC-C3C	-2.33	1.41	1.47
4	B	550	HEM	C2A-C3A	-2.31	1.32	1.38
5	B	601	H4B	C6-N5	2.31	1.51	1.47
5	D	603	H4B	C6-N5	2.28	1.51	1.47
4	B	550	HEM	CAC-C3C	-2.23	1.41	1.47
4	D	550	HEM	C3D-C2D	-2.14	1.32	1.36
4	C	550	HEM	CHD-C1D	-2.02	1.34	1.39
4	C	550	HEM	C1A-NA	-2.01	1.35	1.39

All (37) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	C	602	H4B	C2-N1-C8A	5.09	122.34	113.36
5	D	603	H4B	C2-N1-C8A	5.08	122.33	113.36
5	B	601	H4B	C2-N1-C8A	5.06	122.29	113.36
5	A	600	H4B	C2-N1-C8A	5.00	122.18	113.36
4	B	550	HEM	C3B-C4B-NB	3.66	112.10	109.47
4	C	550	HEM	C3B-C4B-NB	3.65	112.09	109.47
4	D	550	HEM	CMD-C2D-C1D	3.60	130.66	125.03
4	A	550	HEM	C3B-C4B-NB	3.55	112.02	109.47
4	D	550	HEM	CBD-CAD-C3D	-3.24	103.58	112.53
4	D	550	HEM	C3B-C4B-NB	3.23	111.79	109.47
5	B	601	H4B	O10-C10-C9	-3.08	104.68	109.77
5	C	602	H4B	O10-C10-C9	-3.07	104.69	109.77
5	C	602	H4B	N3-C2-N1	-2.79	118.22	123.32
5	B	601	H4B	N3-C2-N1	-2.75	118.29	123.32
4	D	550	HEM	C2A-C1A-NA	2.71	113.16	110.15
5	A	600	H4B	N3-C2-N1	-2.71	118.37	123.32
5	D	603	H4B	N3-C2-N1	-2.70	118.37	123.32

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	600	H4B	O10-C10-C9	-2.69	105.32	109.77
4	C	550	HEM	C2D-C1D-ND	2.69	113.01	109.90
4	B	550	HEM	C4C-C3C-C2C	-2.67	104.50	106.81
5	D	603	H4B	O10-C10-C9	-2.61	105.45	109.77
4	B	550	HEM	C2A-C1A-NA	2.61	113.05	110.15
4	C	550	HEM	C1D-C2D-C3D	-2.61	104.24	106.98
4	C	550	HEM	CMD-C2D-C1D	2.56	129.03	125.03
4	B	550	HEM	C4A-NA-C1A	-2.52	101.72	105.82
4	A	550	HEM	C4D-ND-C1D	-2.49	102.26	105.21
4	B	550	HEM	C2B-C1B-NB	2.47	112.67	109.84
4	C	550	HEM	CAD-C3D-C2D	-2.41	123.35	127.87
4	C	550	HEM	C4B-C3B-C2B	-2.35	105.12	107.28
4	A	550	HEM	CMD-C2D-C1D	2.29	128.61	125.03
4	B	550	HEM	CMA-C3A-C4A	2.23	128.81	125.42
4	C	550	HEM	C4C-NC-C1C	-2.20	102.24	105.82
4	D	550	HEM	CMB-C2B-C1B	2.20	128.47	125.03
4	B	550	HEM	CAA-C2A-C1A	2.19	129.23	124.94
4	C	550	HEM	C4D-ND-C1D	-2.10	102.72	105.21
4	B	550	HEM	C1B-NB-C4B	-2.08	102.74	105.21
4	A	550	HEM	C2D-C1D-ND	2.06	112.29	109.90

There are no chirality outliers.

All (22) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	550	HEM	C2B-C3B-CAB-CBB
4	B	550	HEM	C2B-C3B-CAB-CBB
7	A	880	GOL	O1-C1-C2-C3
7	B	881	GOL	O1-C1-C2-C3
7	A	880	GOL	C1-C2-C3-O3
7	B	881	GOL	C1-C2-C3-O3
7	C	882	GOL	O1-C1-C2-C3
7	D	883	GOL	O1-C1-C2-C3
7	D	883	GOL	C1-C2-C3-O3
7	A	880	GOL	O1-C1-C2-O2
7	A	880	GOL	O2-C2-C3-O3
7	B	881	GOL	O1-C1-C2-O2
7	D	883	GOL	O1-C1-C2-O2
7	D	883	GOL	O2-C2-C3-O3
4	A	550	HEM	C4B-C3B-CAB-CBB
4	B	550	HEM	C4B-C3B-CAB-CBB
4	D	550	HEM	C4B-C3B-CAB-CBB

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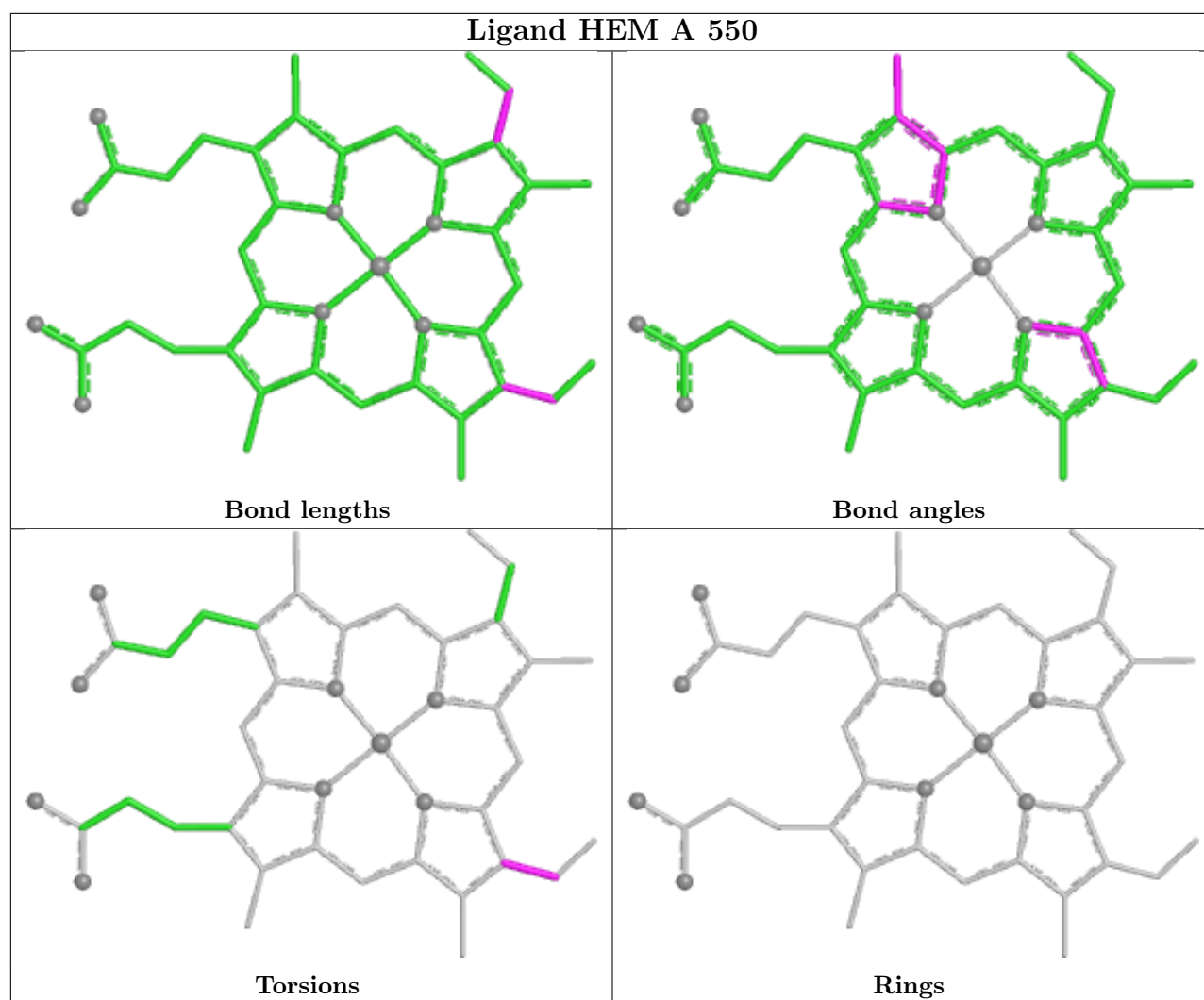
Mol	Chain	Res	Type	Atoms
7	B	881	GOL	O2-C2-C3-O3
4	B	550	HEM	CAA-CBA-CGA-O1A
4	B	550	HEM	CAA-CBA-CGA-O2A
4	C	550	HEM	CAA-CBA-CGA-O2A
4	C	550	HEM	CAA-CBA-CGA-O1A

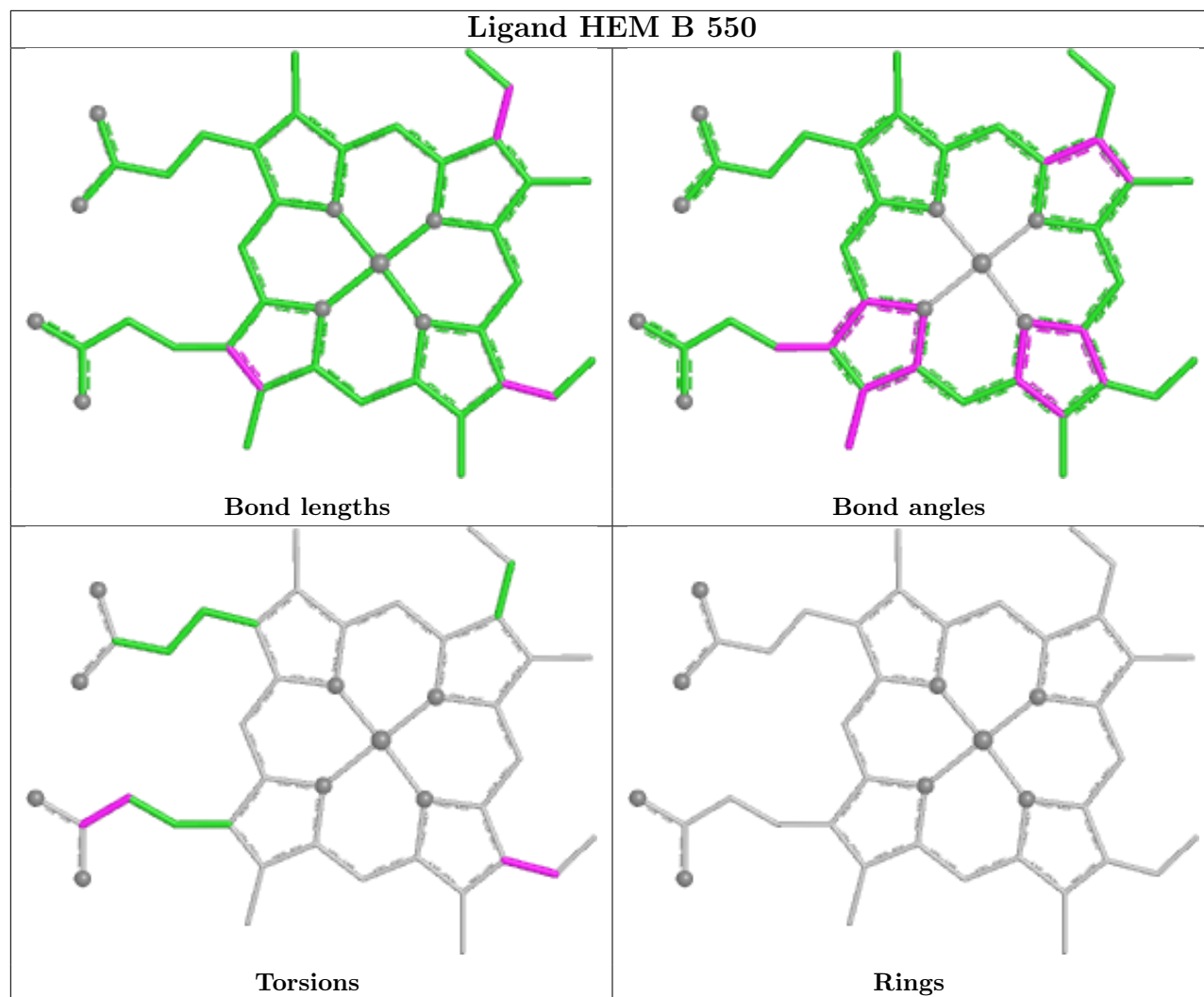
There are no ring outliers.

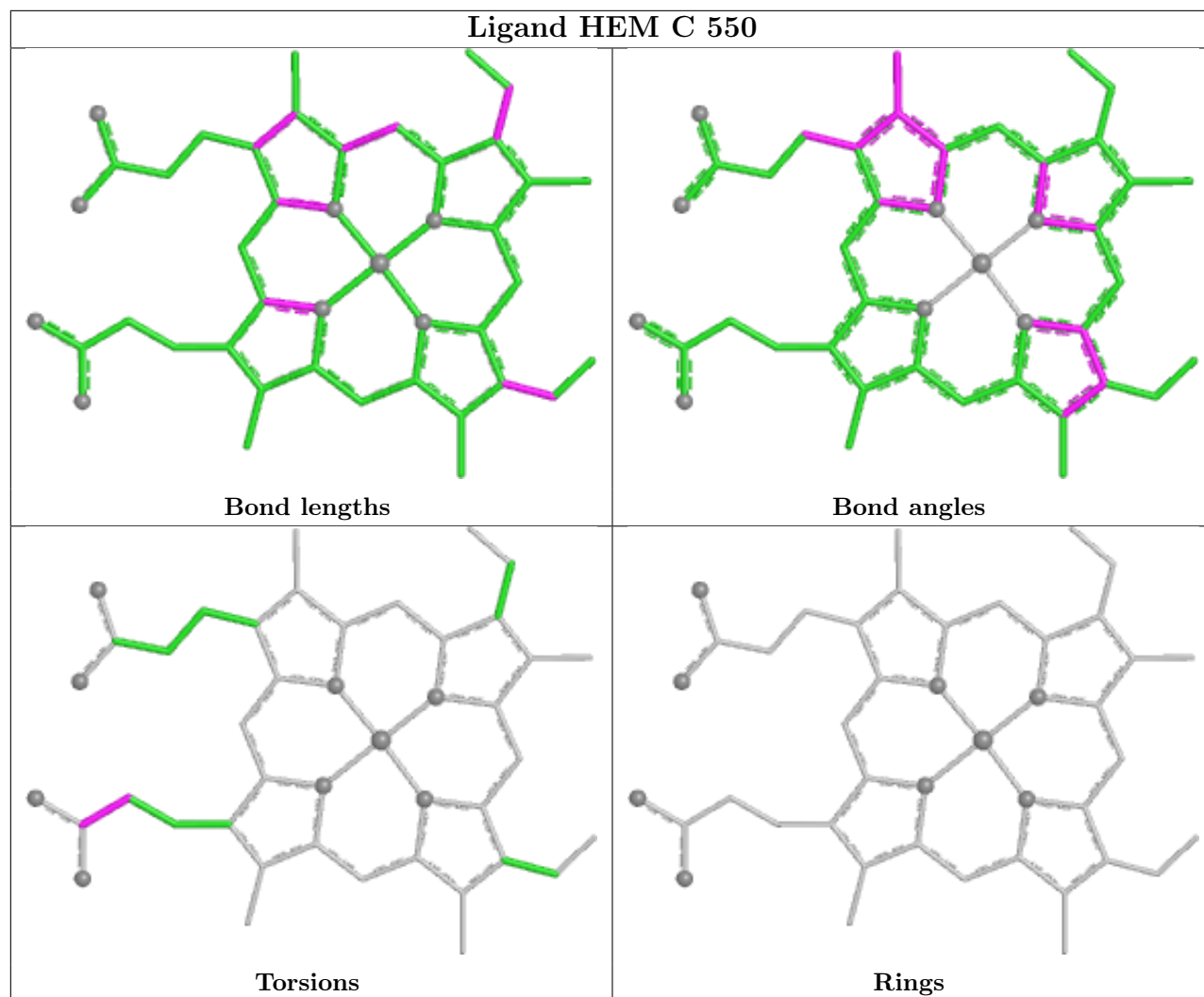
6 monomers are involved in 11 short contacts:

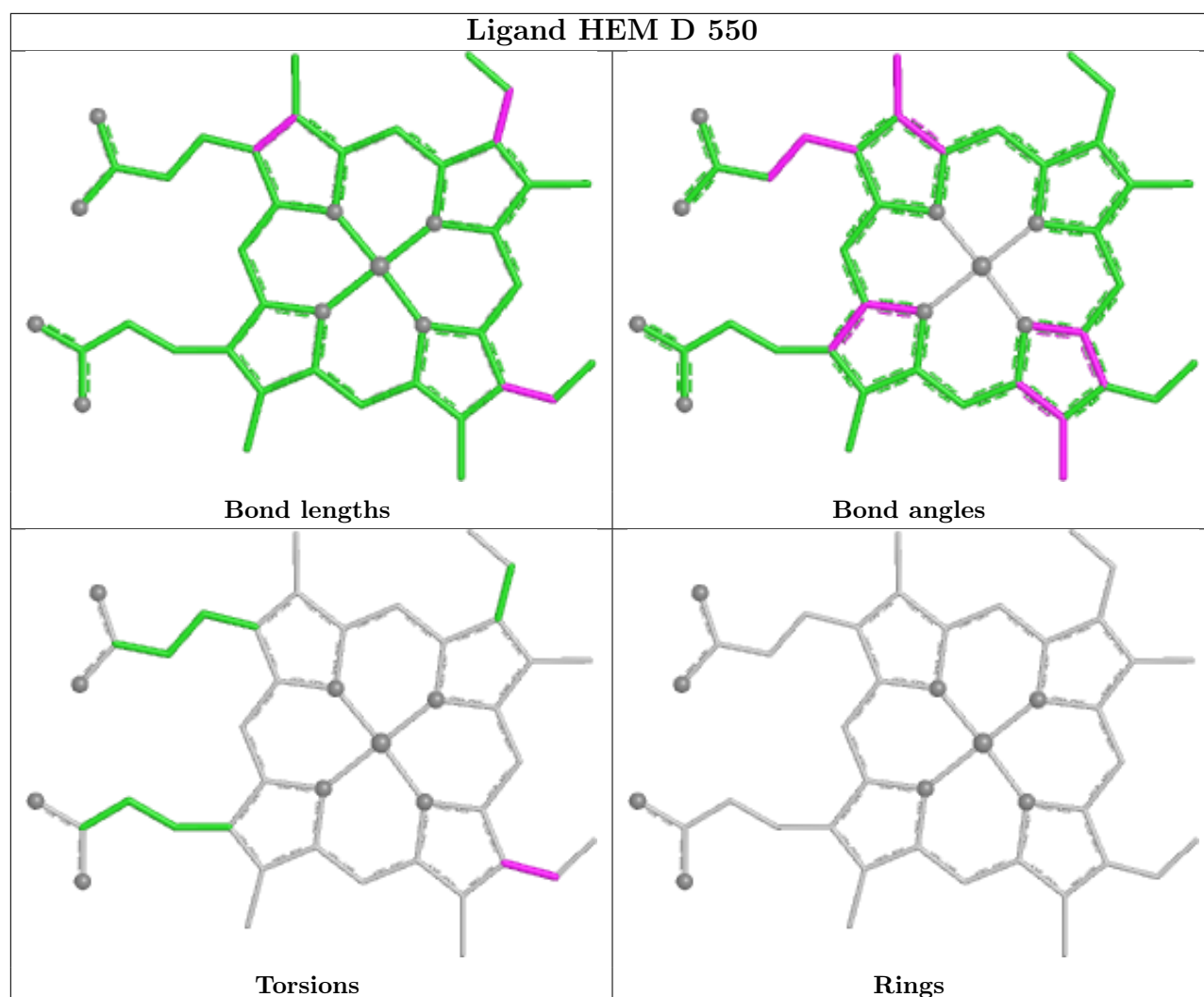
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	550	HEM	1	0
4	B	550	HEM	2	0
7	D	883	GOL	3	0
2	D	923	SO4	1	0
5	C	602	H4B	1	0
4	D	550	HEM	3	0

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.