



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 5, 2026 – 01:01 PM EDT

PDB ID : 1LGA / pdb_00001lga
Title : CRYSTALLOGRAPHIC REFINEMENT OF LIGNIN PEROXIDASE AT 2
ANGSTROMS
Authors : Poulos, T.L.; Edwards, S.L.; Wariishi, H.; Gold, M.H.
Deposited on : 1992-12-08
Resolution : 2.03 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

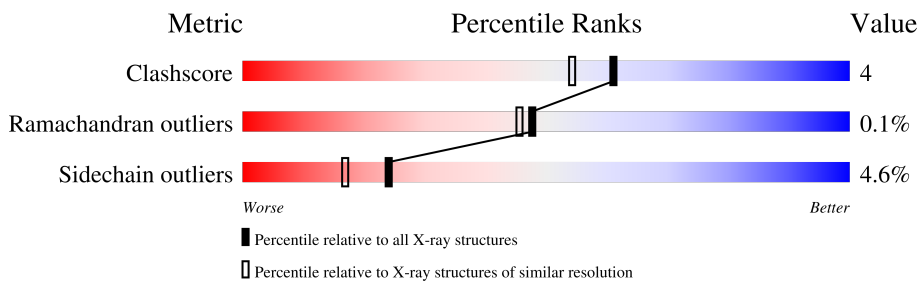
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.03 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	1022 (2.02-2.02)
Ramachandran outliers	187476	1014 (2.02-2.02)
Sidechain outliers	187428	1014 (2.02-2.02)

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	343	
1	B	343	

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 5693 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 LIGNIN PEROXIDASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	343	Total	C	N	O	S	0	0	0
			2556	1615	428	497	16			
1	B	341	Total	C	N	O	S	0	0	0
			2545	1609	426	494	16			

- Molecule 2 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).

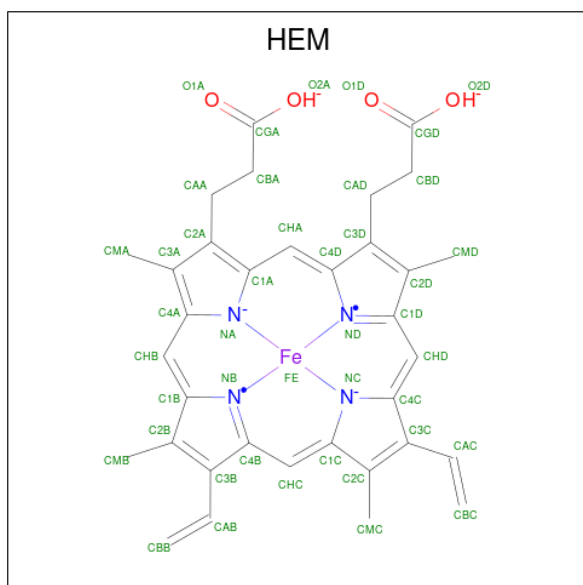


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	B	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 3 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	2	Total Ca 2 2	0	0
3	B	2	Total Ca 2 2	0	0

- Molecule 4 is PROTOPORPHYRIN IX CONTAINING FE (CCD ID: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).

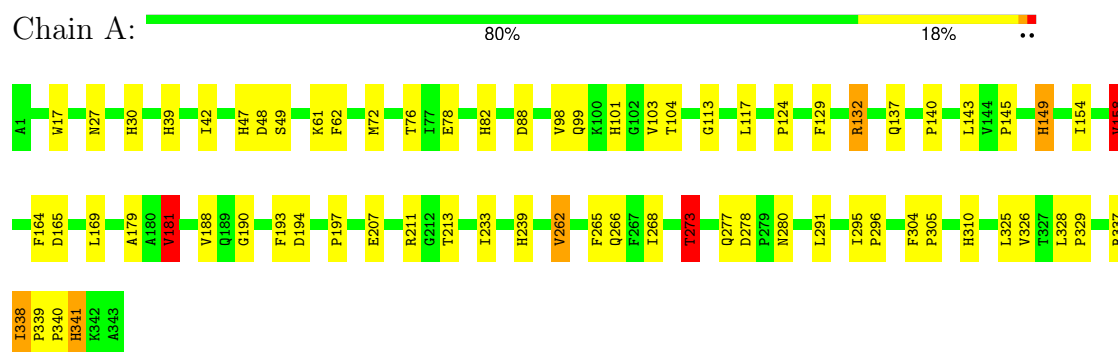


3 Residue-property plots [i](#)

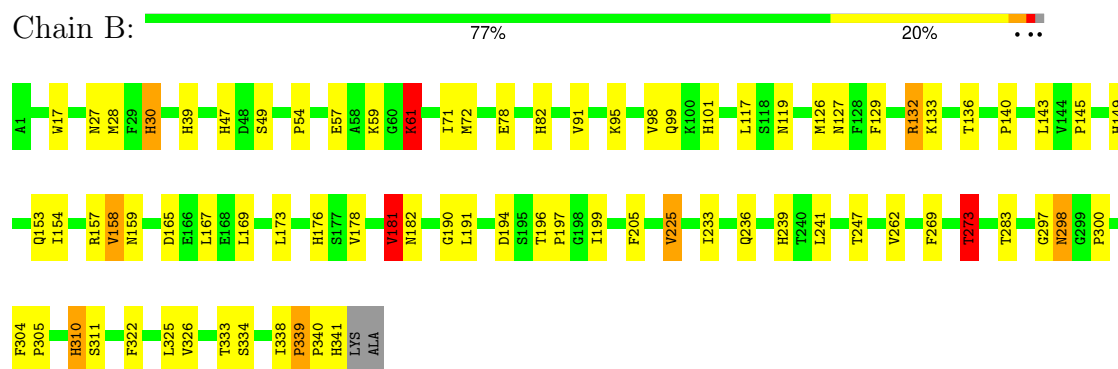
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

Note EDS was not executed.

• Molecule 1: LIGNIN PEROXIDASE



• Molecule 1: LIGNIN PEROXIDASE



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	44.70Å 77.50Å 100.00Å 90.00° 101.00° 90.00°	Depositor
Resolution (Å)	(Not available) – 2.03	Depositor
% Data completeness (in resolution range)	(Not available) ((Not available)-2.03)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR	Depositor
R, R_{free}	0.150 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	5693	wwPDB-VP
Average B, all atoms (Å ²)	13.0	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: CA, NAG, 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	1.07	13/2626 (0.5%)	1.72	46/3578 (1.3%)
1	B	1.11	13/2615 (0.5%)	1.73	50/3564 (1.4%)
All	All	1.09	26/5241 (0.5%)	1.73	96/7142 (1.3%)

All (26) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	47	HIS	CD2-NE2	-7.43	1.29	1.37
1	B	239	HIS	CD2-NE2	-7.21	1.29	1.37
1	A	47	HIS	CD2-NE2	-6.98	1.30	1.37
1	A	39	HIS	CD2-NE2	-6.86	1.30	1.37
1	B	101	HIS	CD2-NE2	-6.83	1.30	1.37
1	B	341	HIS	CD2-NE2	-6.79	1.30	1.37
1	B	310	HIS	CD2-NE2	-6.72	1.30	1.37
1	A	310	HIS	CD2-NE2	-6.60	1.30	1.37
1	A	101	HIS	CD2-NE2	-6.50	1.30	1.37
1	B	149	HIS	CD2-NE2	-6.47	1.30	1.37
1	B	181	VAL	CA-CB	6.43	1.62	1.54
1	B	30	HIS	CD2-NE2	-6.39	1.30	1.37
1	A	82	HIS	CD2-NE2	-6.33	1.30	1.37
1	B	39	HIS	CD2-NE2	-6.21	1.31	1.37
1	B	326	VAL	CA-CB	5.99	1.61	1.54
1	A	341	HIS	CD2-NE2	-5.96	1.31	1.37
1	B	71	ILE	CA-CB	5.82	1.62	1.54
1	A	239	HIS	CD2-NE2	-5.80	1.31	1.37
1	A	149	HIS	CD2-NE2	-5.72	1.31	1.37
1	A	30	HIS	CD2-NE2	-5.59	1.31	1.37
1	B	82	HIS	CD2-NE2	-5.50	1.31	1.37
1	A	188	VAL	CA-CB	5.45	1.61	1.55
1	A	181	VAL	CA-CB	5.36	1.61	1.54
1	A	101	HIS	CG-ND1	-5.27	1.32	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	178	VAL	CA-CB	5.20	1.61	1.54
1	A	137	GLN	CD-NE2	-5.13	1.22	1.33

All (96) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	132	ARG	NE-CZ-NH2	-12.81	107.67	119.20
1	B	225	VAL	N-CA-CB	-11.52	99.55	112.45
1	B	194	ASP	CA-CB-CG	8.99	121.59	112.60
1	A	194	ASP	CA-CB-CG	8.95	121.55	112.60
1	A	132	ARG	NE-CZ-NH2	-8.89	111.20	119.20
1	B	298	ASN	N-CA-C	-8.68	96.97	110.42
1	B	132	ARG	NE-CZ-NH1	8.46	129.96	121.50
1	A	99	GLN	OE1-CD-NE2	-8.21	114.39	122.60
1	A	338	ILE	N-CA-C	-8.18	98.85	107.60
1	B	298	ASN	CA-CB-CG	7.83	120.43	112.60
1	A	296	PRO	N-CA-C	7.56	123.47	113.57
1	A	273	THR	N-CA-CB	-7.51	97.55	110.32
1	A	132	ARG	NE-CZ-NH1	7.50	129.00	121.50
1	B	159	ASN	CA-CB-CG	-7.47	105.13	112.60
1	A	49	SER	N-CA-C	7.42	119.01	111.07
1	B	98	VAL	N-CA-C	-7.40	103.32	110.42
1	B	262	VAL	N-CA-CB	-7.33	101.19	110.57
1	B	132	ARG	CB-CG-CD	7.25	127.98	111.30
1	B	127	ASN	OD1-CG-ND2	-7.09	115.51	122.60
1	B	158	VAL	N-CA-CB	-7.05	100.00	110.58
1	A	78	GLU	N-CA-C	7.05	119.04	111.36
1	B	165	ASP	CA-CB-CG	6.96	119.56	112.60
1	B	61	LYS	N-CA-C	6.74	119.47	108.55
1	A	266	GLN	OE1-CD-NE2	-6.60	116.00	122.60
1	A	158	VAL	N-CA-CB	-6.60	99.37	110.65
1	B	101	HIS	CA-CB-CG	-6.53	107.27	113.80
1	B	273	THR	N-CA-CB	-6.52	100.18	110.22
1	B	225	VAL	CB-CA-C	6.50	120.44	111.34
1	A	76	THR	N-CA-C	-6.48	104.22	111.28
1	B	341	HIS	CB-CG-CD2	-6.47	122.80	131.20
1	A	304	PHE	N-CA-C	-6.39	100.35	109.62
1	A	165	ASP	CA-CB-CG	6.35	118.95	112.60
1	A	273	THR	OG1-CB-CG2	6.32	121.94	109.30
1	A	101	HIS	CB-CG-CD2	-6.30	123.01	131.20
1	A	280	ASN	OD1-CG-ND2	-6.22	116.38	122.60
1	B	334	SER	O-C-N	-6.20	116.06	123.31

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	88	ASP	CA-CB-CG	6.17	118.77	112.60
1	B	239	HIS	CB-CG-CD2	-6.15	123.20	131.20
1	A	132	ARG	CB-CG-CD	6.15	125.44	111.30
1	B	196	THR	CA-C-N	6.09	125.77	119.56
1	B	196	THR	C-N-CA	6.09	125.77	119.56
1	A	48	ASP	CA-C-N	6.07	128.33	120.44
1	A	48	ASP	C-N-CA	6.07	128.33	120.44
1	A	103	VAL	O-C-N	-6.06	116.48	122.12
1	A	30	HIS	CB-CG-CD2	-5.90	123.53	131.20
1	B	132	ARG	CG-CD-NE	-5.90	99.02	112.00
1	B	28	MET	N-CA-C	-5.87	106.03	113.43
1	B	236	GLN	OE1-CD-NE2	-5.86	116.74	122.60
1	A	278	ASP	CA-CB-CG	5.78	118.38	112.60
1	B	47	HIS	CB-CG-CD2	-5.75	123.72	131.20
1	B	30	HIS	CB-CG-CD2	-5.74	123.74	131.20
1	B	17	TRP	CG-CD2-CE3	5.68	139.59	133.90
1	A	27	ASN	OD1-CG-ND2	-5.68	116.92	122.60
1	B	305	PRO	O-C-N	-5.66	118.71	121.31
1	A	132	ARG	CG-CD-NE	-5.64	99.60	112.00
1	A	295	ILE	N-CA-C	-5.60	103.46	108.95
1	B	49	SER	N-CA-C	5.59	117.18	111.14
1	B	305	PRO	CA-C-O	-5.56	116.90	120.90
1	B	297	GLY	CA-C-O	5.54	126.90	122.52
1	A	99	GLN	CG-CD-NE2	5.54	124.70	116.40
1	B	191	LEU	CA-C-O	-5.51	114.85	119.76
1	B	78	GLU	N-CA-C	5.50	117.35	111.36
1	B	304	PHE	CA-C-N	5.49	123.67	119.66
1	B	304	PHE	C-N-CA	5.49	123.67	119.66
1	A	47	HIS	CB-CG-CD2	-5.47	124.09	131.20
1	B	119	ASN	OD1-CG-ND2	-5.47	117.13	122.60
1	A	104	THR	CA-C-N	5.45	125.53	119.32
1	A	104	THR	C-N-CA	5.45	125.53	119.32
1	B	269	PHE	N-CA-C	-5.39	105.32	111.14
1	B	132	ARG	CD-NE-CZ	5.37	131.91	124.40
1	A	326	VAL	CB-CA-C	-5.34	105.47	111.45
1	A	117	LEU	CA-C-O	-5.33	114.77	120.42
1	B	133	LYS	CB-CG-CD	-5.32	99.06	111.30
1	B	27	ASN	CA-CB-CG	-5.32	107.28	112.60
1	B	205	PHE	CA-CB-CG	-5.30	108.50	113.80
1	A	278	ASP	CA-C-N	5.30	125.11	119.28
1	A	278	ASP	C-N-CA	5.30	125.11	119.28
1	B	197	PRO	N-CA-C	5.26	120.68	113.84

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	341	HIS	CB-CG-CD2	-5.25	124.37	131.20
1	A	113	GLY	O-C-N	-5.24	117.11	122.19
1	A	17	TRP	CG-CD2-CE3	5.22	139.12	133.90
1	B	339	PRO	N-CA-C	5.20	117.05	110.70
1	B	101	HIS	CB-CG-CD2	-5.18	124.46	131.20
1	A	42	ILE	N-CA-C	-5.16	105.68	110.53
1	B	300	PRO	CB-CA-C	5.14	117.72	110.98
1	B	176	HIS	CG-CD2-NE2	-5.13	102.07	107.20
1	A	262	VAL	N-CA-CB	-5.10	104.04	110.57
1	B	17	TRP	CB-CG-CD1	-5.10	119.25	126.90
1	B	158	VAL	CB-CA-C	5.08	119.23	112.22
1	A	239	HIS	CB-CG-CD2	-5.07	124.61	131.20
1	B	95	LYS	CA-C-O	-5.07	113.68	118.44
1	A	17	TRP	CE2-CD2-CG	-5.06	101.13	107.20
1	A	310	HIS	CB-CG-CD2	-5.06	124.63	131.20
1	A	305	PRO	CA-C-N	5.04	125.32	119.93
1	A	305	PRO	C-N-CA	5.04	125.32	119.93
1	A	207	GLU	N-CA-C	5.01	119.10	112.89

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	2556	0	2411	22	0
1	B	2545	0	2405	20	0
2	A	14	0	13	0	0
2	B	14	0	13	0	0
3	A	2	0	0	0	0
3	B	2	0	0	0	0
4	A	43	0	30	1	0
4	B	43	0	30	1	0
5	A	230	0	0	1	0
5	B	244	0	0	2	0
All	All	5693	0	4902	42	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 4.

All (42) 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:181:VAL:HG21	1:A:233:ILE:HD13	1.75	0.68
1:A:158:VAL:HG13	1:A:164:PHE:HB2	1.86	0.58
1:A:265:PHE:HA	1:A:268:ILE:HG22	1.88	0.56
1:B:132:ARG:NH1	1:B:273:THR:HG23	2.25	0.51
1:B:145:PRO:HB3	1:B:154:ILE:HD12	1.92	0.51
1:A:140:PRO:HG2	1:A:143:LEU:HD11	1.93	0.51
1:B:310:HIS:HE1	1:B:322:PHE:O	1.94	0.50
1:A:145:PRO:HB3	1:A:154:ILE:HD12	1.93	0.49
1:B:273:THR:HG21	5:B:473:HOH:O	2.13	0.49
1:B:129:PHE:O	1:B:132:ARG:NH2	2.45	0.49
1:B:181:VAL:HG21	1:B:233:ILE:HD13	1.95	0.49
1:B:182:ASN:ND2	1:B:340:PRO:HB3	2.28	0.48
1:A:149:HIS:HB2	1:A:154:ILE:HD11	1.94	0.48
1:A:181:VAL:CG2	1:A:233:ILE:HD13	2.44	0.48
1:B:72:MET:HG3	1:B:91:VAL:HG13	1.95	0.48
1:A:190:GLY:HA3	1:A:338:ILE:O	2.15	0.46
1:A:154:ILE:HG21	4:A:396:HEM:HBB1	1.97	0.46
1:B:117:LEU:HD13	1:B:126:MET:HE2	1.97	0.46
1:B:190:GLY:HA3	1:B:338:ILE:O	2.17	0.45
1:A:129:PHE:O	1:A:132:ARG:NH2	2.50	0.44
1:B:339:PRO:HA	1:B:340:PRO:HD2	1.82	0.44
1:B:241:LEU:HD22	1:B:247:THR:HG21	1.99	0.44
1:B:140:PRO:HG2	1:B:143:LEU:HD11	2.00	0.43
1:A:179:ALA:HB3	1:A:193:PHE:HD2	1.84	0.43
1:A:197:PRO:HG3	1:A:338:ILE:HD11	2.01	0.43
1:A:211:ARG:NH1	1:A:213:THR:HG22	2.34	0.43
1:A:339:PRO:HA	1:A:340:PRO:HD2	1.85	0.42
1:A:72:MET:SD	1:A:98:VAL:HG21	2.59	0.42
1:B:199:ILE:HG21	1:B:199:ILE:HD13	1.87	0.42
1:B:338:ILE:HA	1:B:339:PRO:HD2	1.85	0.42
1:A:132:ARG:HG2	1:A:277:GLN:CD	2.45	0.42
1:A:145:PRO:HB3	1:A:154:ILE:CD1	2.50	0.41
1:B:173:LEU:HD23	1:B:173:LEU:HA	1.92	0.41
1:A:124:PRO:HB3	1:A:262:VAL:HG13	2.01	0.41
1:B:129:PHE:HA	1:B:283:THR:O	2.20	0.41
1:B:57:GLU:HA	1:B:61:LYS:O	2.20	0.41
1:A:273:THR:HG21	5:A:485:HOH:O	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:167:LEU:HD23	5:B:439:HOH:O	2.21	0.41
1:B:181:VAL:HB	4:B:396:HEM:HAA2	2.03	0.40
1:A:268:ILE:HD12	1:A:268:ILE:HA	1.93	0.40
1:A:328:LEU:HA	1:A:329:PRO:HD3	1.84	0.40
1:A:61:LYS:H	1:A:61:LYS:HD3	1.85	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	341/343 (99%)	331 (97%)	9 (3%)	1 (0%)	36	31
1	B	339/343 (99%)	331 (98%)	8 (2%)	0	100	100
All	All	680/686 (99%)	662 (97%)	17 (2%)	1 (0%)	48	45

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	62	PHE

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.

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	273/274 (100%)	265 (97%)	8 (3%)	37	34
1	B	273/274 (100%)	256 (94%)	17 (6%)	16	9
All	All	546/548 (100%)	521 (95%)	25 (5%)	24	17

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	158	VAL
1	A	169	LEU
1	A	181	VAL
1	A	273	THR
1	A	291	LEU
1	A	325	LEU
1	A	337	ARG
1	A	341	HIS
1	B	30	HIS
1	B	54	PRO
1	B	59	LYS
1	B	61	LYS
1	B	99	GLN
1	B	136	THR
1	B	153	GLN
1	B	157	ARG
1	B	158	VAL
1	B	169	LEU
1	B	181	VAL
1	B	225	VAL
1	B	273	THR
1	B	298	ASN
1	B	311	SER
1	B	325	LEU
1	B	333	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	27	ASN
1	A	94	GLN

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Mol	Chain	Res	Type
1	A	99	GLN
1	A	127	ASN
1	A	310	HIS
1	B	33	GLN
1	B	94	GLN
1	B	277	GLN
1	B	310	HIS

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 ⓘ

Of 8 ligands modelled in this entry, 4 are monoatomic - leaving 4 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
4	HEM	B	396	5,1	50,50,50	1.31	5 (10%)	67,82,82	1.36	10 (14%)
4	HEM	A	396	5,1	50,50,50	1.16	6 (12%)	67,82,82	1.04	4 (5%)
2	NAG	B	397	1	14,14,15	0.94	1 (7%)	17,19,21	1.20	3 (17%)
2	NAG	A	397	1	14,14,15	0.93	0	17,19,21	1.13	1 (5%)

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
4	HEM	B	396	5,1	-	2/14/54/54	-
4	HEM	A	396	5,1	-	3/14/54/54	-
2	NAG	B	397	1	-	0/6/23/26	0/1/1/1
2	NAG	A	397	1	-	0/6/23/26	0/1/1/1

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	396	HEM	CBC-CAC	3.22	1.45	1.30
4	B	396	HEM	C4D-ND	-3.16	1.34	1.40
4	B	396	HEM	CAB-C3B	-2.94	1.39	1.47
4	A	396	HEM	CBB-CAB	2.89	1.44	1.30
4	B	396	HEM	CAC-C3C	-2.61	1.40	1.47
4	A	396	HEM	CAC-C3C	-2.55	1.40	1.47
4	A	396	HEM	CBC-CAC	2.55	1.42	1.30
4	B	396	HEM	CBB-CAB	2.54	1.42	1.30
4	A	396	HEM	CAB-C3B	-2.48	1.40	1.47
4	A	396	HEM	O2D-CGD	-2.44	1.22	1.30
2	B	397	NAG	C3-C2	-2.25	1.47	1.52
4	A	396	HEM	O2A-CGA	-2.14	1.23	1.30

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	396	HEM	C3B-C4B-NB	4.52	112.71	109.47
4	A	396	HEM	C3B-C4B-NB	3.97	112.32	109.47
4	B	396	HEM	C1D-C2D-C3D	-3.58	103.22	106.98
2	A	397	NAG	C1-C2-N2	-3.29	105.25	110.43
4	B	396	HEM	CAC-C3C-C4C	2.90	131.74	124.82
4	A	396	HEM	C4C-C3C-C2C	-2.72	104.45	106.81
2	B	397	NAG	C2-N2-C7	-2.70	119.28	122.90
2	B	397	NAG	O4-C4-C3	-2.60	104.24	110.38
4	A	396	HEM	O2D-CGD-CBD	2.43	121.68	114.00
4	B	396	HEM	CMD-C2D-C1D	2.32	128.66	125.03
2	B	397	NAG	C1-C2-N2	-2.27	106.86	110.43
4	B	396	HEM	C2B-C1B-NB	2.23	112.41	109.84
4	B	396	HEM	O1D-CGD-CBD	-2.17	116.21	123.09
4	A	396	HEM	C1D-C2D-C3D	-2.16	104.71	106.98
4	B	396	HEM	O2D-CGD-CBD	2.08	120.56	114.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	396	HEM	C4C-C3C-C2C	-2.06	105.03	106.81
4	B	396	HEM	C2D-C1D-ND	2.05	112.27	109.90
4	B	396	HEM	C4D-C3D-C2D	2.01	109.82	106.89

There are no chirality outliers.

All (5) torsion outliers are listed below:

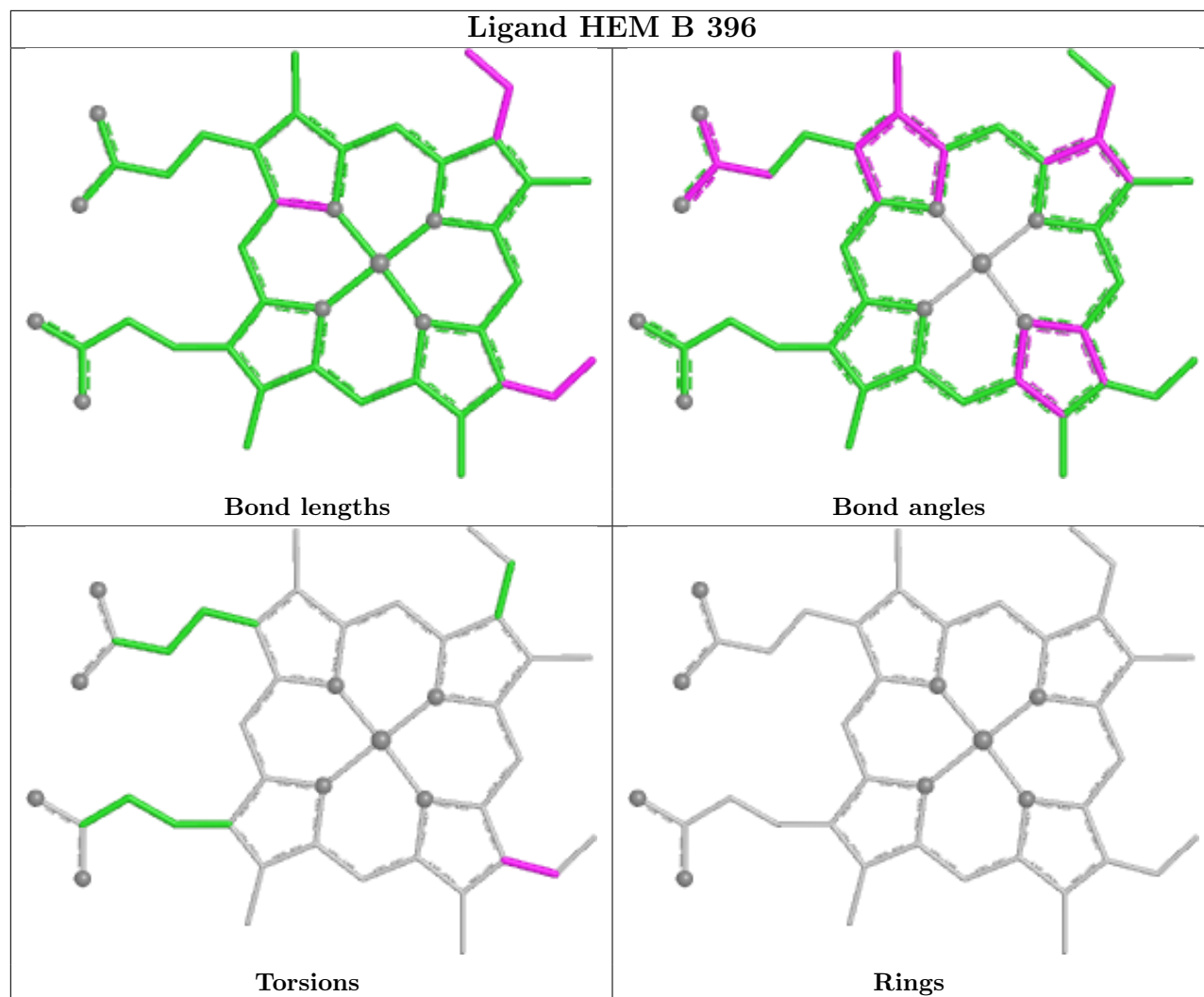
Mol	Chain	Res	Type	Atoms
4	A	396	HEM	C2C-C3C-CAC-CBC
4	A	396	HEM	C2B-C3B-CAB-CBB
4	B	396	HEM	C2B-C3B-CAB-CBB
4	B	396	HEM	C4B-C3B-CAB-CBB
4	A	396	HEM	C4B-C3B-CAB-CBB

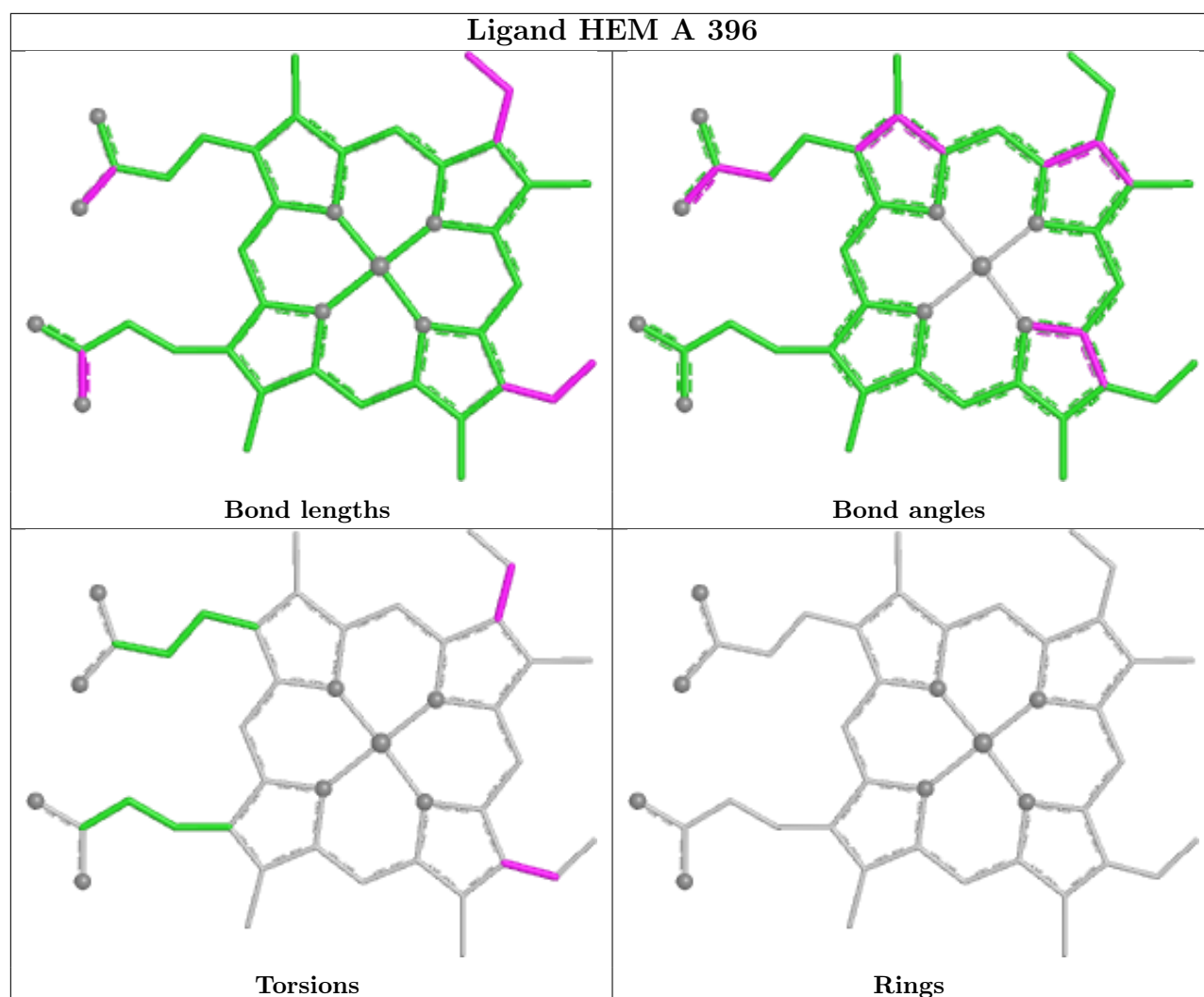
There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	396	HEM	1	0
4	A	396	HEM	1	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.