



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 18, 2026 – 07:33 AM UTC

PDB ID : 5JPR / pdb_00005jpr
Title : Neutron Structure of Compound II of Ascorbate Peroxidase
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Deposited on : 2016-05-04
Resolution : 1.81 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)	:	2.0
EDS	:	FAILED
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

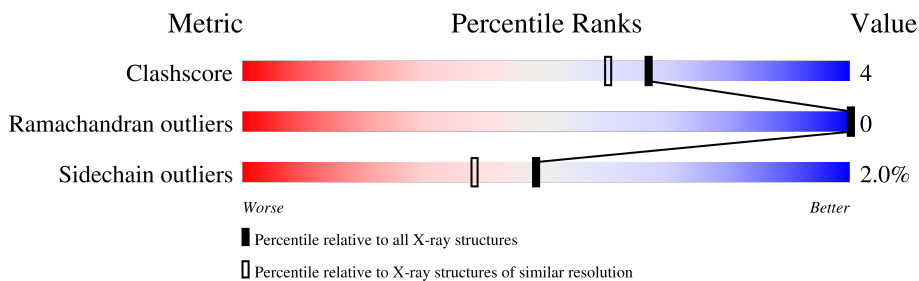
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION, NEUTRON DIFFRACTION

The reported resolution of this entry is 1.81 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	190562	8479 (1.80-1.80)
Ramachandran outliers	187476	8391 (1.80-1.80)
Sidechain outliers	187428	8390 (1.80-1.80)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$

Note EDS failed to run properly.

Mol	Chain	Length	Quality of chain
1	A	261	

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
4	SO4	A	303	-	-	X	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 5180 atoms, of which 1893 are hydrogens and 944 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.

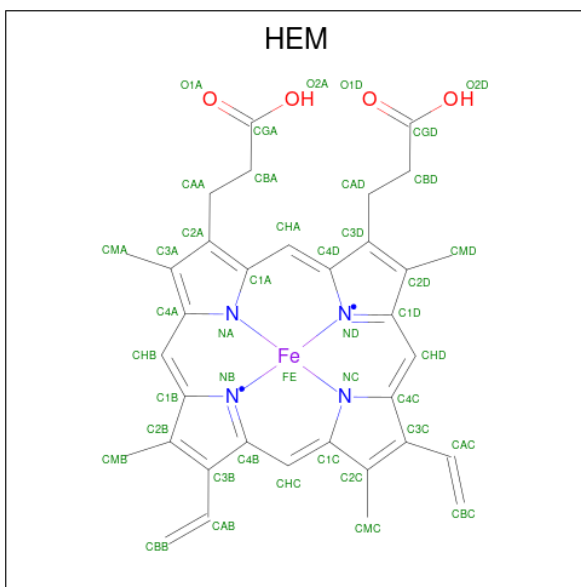
- Molecule 1 is a protein called Ascorbate peroxidase.

Mol	Chain	Residues	Atoms							ZeroOcc	AltConf	Trace
1	A	248	Total	C	D	H	N	O	S	141	228	0
			4158	1214	396	1863	321	361	3			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-10	MET	-	initiating methionine	UNP Q43758
A	-9	ARG	-	expression tag	UNP Q43758
A	-8	GLY	-	expression tag	UNP Q43758
A	-7	SER	-	expression tag	UNP Q43758
A	-6	HIS	-	expression tag	UNP Q43758
A	-5	HIS	-	expression tag	UNP Q43758
A	-4	HIS	-	expression tag	UNP Q43758
A	-3	HIS	-	expression tag	UNP Q43758
A	-2	HIS	-	expression tag	UNP Q43758
A	-1	HIS	-	expression tag	UNP Q43758
A	0	GLY	-	expression tag	UNP Q43758
A	1	SER	-	expression tag	UNP Q43758

- Molecule 2 is PROTOPORPHYRIN IX CONTAINING FE (CCD ID: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).

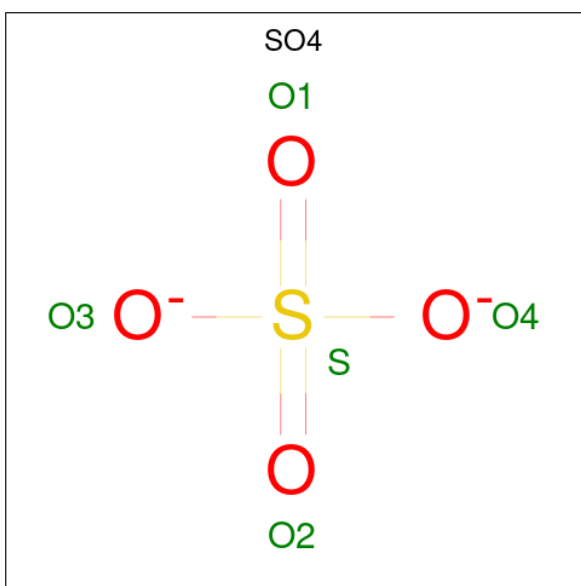


Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total	C	Fe	H	N	O	0	0
			73	34	1	30	4	4		

- Molecule 3 is POTASSIUM ION (CCD ID: K) (formula: K).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	K	0	0
			1	1		

- Molecule 4 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

- Molecule 5 is water.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	390	Total	D	O	0	0
			938	548	390		

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 failed to run properly.

- Molecule 1: Ascorbate peroxidase

Chain A:  90% 5% 5%



4 Data and refinement statistics

EDS failed to run properly - this section is therefore incomplete.

Property	Value	Source
Space group	P 42 21 2	Depositor
Cell constants a, b, c, α , β , γ	82.09Å 82.09Å 75.16Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.91 – 1.81	Depositor
% Data completeness (in resolution range)	98.7 (19.91-1.81)	Depositor
R_{merge}	0.18	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.96 (at 1.80Å)	Xtriage
Refinement program	PHENIX 1.7.3_928	Depositor
R, R_{free}	0.155 , 0.216	Depositor
Wilson B-factor (Å ²)	13.6	Xtriage
Anisotropy	0.097	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	5180	wwPDB-VP
Average B, all atoms (Å ²)	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.82% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: DOD, K, SO4, 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	1.38	6/3744 (0.2%)	1.25	21/5070 (0.4%)

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	210[A]	ALA	CA-CB	-7.36	1.40	1.53
1	A	210[B]	ALA	CA-CB	-7.36	1.40	1.53
1	A	79[A]	ARG	CZ-NH1	5.96	1.41	1.32
1	A	79[B]	ARG	CZ-NH1	5.96	1.41	1.32
1	A	86[A]	ALA	CA-C	-5.05	1.45	1.52
1	A	86[B]	ALA	CA-C	-5.05	1.45	1.52

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	110[A]	GLY	N-CA-C	6.73	120.37	112.23
1	A	110[B]	GLY	N-CA-C	6.73	120.37	112.23
1	A	182[A]	ASN	CA-C-N	6.68	127.53	120.12
1	A	182[A]	ASN	C-N-CA	6.68	127.53	120.12
1	A	182[B]	ASN	CA-C-N	6.68	127.53	120.12
1	A	182[B]	ASN	C-N-CA	6.68	127.53	120.12
1	A	205[A]	LEU	CA-C-N	-6.47	112.16	119.28
1	A	205[A]	LEU	C-N-CA	-6.47	112.16	119.28
1	A	205[B]	LEU	CA-C-N	-6.47	112.16	119.28
1	A	205[B]	LEU	C-N-CA	-6.47	112.16	119.28
1	A	195[A]	LEU	N-CA-C	6.41	119.08	111.71
1	A	195[B]	LEU	N-CA-C	6.41	119.08	111.71
1	A	79[A]	ARG	NE-CZ-NH2	-6.02	113.78	119.20
1	A	79[B]	ARG	NE-CZ-NH2	-6.02	113.78	119.20
1	A	125	PRO	O-C-N	5.98	124.06	121.31

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	33[A]	ALA	CA-C-N	-5.86	112.54	119.05
1	A	33[A]	ALA	C-N-CA	-5.86	112.54	119.05
1	A	33[B]	ALA	CA-C-N	-5.86	112.54	119.05
1	A	33[B]	ALA	C-N-CA	-5.86	112.54	119.05
1	A	3[A]	LYS	N-CA-C	5.76	119.21	109.76
1	A	3[B]	LYS	N-CA-C	5.76	119.21	109.76

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	2295	1863	135	0	0
2	A	43	30	30	0	0
3	A	1	0	0	0	0
4	A	10	0	0	3	0
5	A	938	0	0	9	2
All	All	3287	1893	165	9	2

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 (9) 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)
4:A:303:SO4:O3	5:A:402:DOD:O	1.97	0.82
4:A:304:SO4:O2	5:A:403:DOD:O	2.21	0.50
4:A:303:SO4:O2	5:A:404:DOD:O	2.25	0.43

All (2) 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)
5:A:421:DOD:O	5:A:421:DOD:O[2_565]	1.54	0.66
5:A:628:DOD:O	5:A:667:DOD:O[6_565]	2.04	0.16

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	473/261 (181%)	467 (99%)	6 (1%)	0	100	100

There are no Ramachandran outliers to report.

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	373/206 (181%)	365 (98%)	8 (2%)	47	36

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	41[A]	TRP
1	A	41[B]	TRP
1	A	61[A]	LYS
1	A	61[B]	LYS
1	A	84[A]	LEU
1	A	84[B]	LEU
1	A	200[A]	GLU

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Mol	Chain	Res	Type
1	A	200[B]	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 4 ligands modelled in this entry, 1 is monoatomic - leaving 3 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	HEM	A	301	5,1	50,50,50	1.79	12 (24%)	67,82,82	1.81	16 (23%)
4	SO4	A	303	-	4,4,4	0.55	0	6,6,6	0.72	0
4	SO4	A	304	-	4,4,4	0.86	0	6,6,6	1.22	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
2	HEM	A	301	5,1	-	3/14/54/54	-

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	301	HEM	C3D-C2D	6.53	1.50	1.36
2	A	301	HEM	FE-ND	3.69	2.06	1.94
2	A	301	HEM	CHD-C1D	-3.48	1.31	1.39
2	A	301	HEM	C1C-NC	-3.19	1.33	1.39
2	A	301	HEM	C3B-C4B	2.70	1.50	1.44
2	A	301	HEM	C1A-NA	-2.46	1.35	1.39
2	A	301	HEM	CMA-C3A	2.41	1.55	1.50
2	A	301	HEM	C2A-C3A	-2.34	1.32	1.38
2	A	301	HEM	CHC-C4B	-2.14	1.34	1.39
2	A	301	HEM	CMD-C2D	2.11	1.55	1.50
2	A	301	HEM	FE-NB	2.07	2.01	1.94
2	A	301	HEM	CAC-C3C	2.03	1.52	1.47

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	301	HEM	C1B-NB-C4B	4.84	110.94	105.21
2	A	301	HEM	CMD-C2D-C1D	4.62	132.25	125.03
2	A	301	HEM	C3B-C4B-NB	-4.40	106.31	109.47
2	A	301	HEM	CHB-C4A-NA	4.09	131.27	123.86
2	A	301	HEM	C4D-ND-C1D	3.68	109.56	105.21
2	A	301	HEM	CHA-C4D-ND	3.49	128.68	124.37
2	A	301	HEM	C2A-C1A-NA	-2.80	107.05	110.15
2	A	301	HEM	O2A-CGA-CBA	2.45	121.75	114.00
2	A	301	HEM	C1A-CHA-C4D	-2.39	120.62	126.25
2	A	301	HEM	CHA-C4D-C3D	-2.27	121.05	125.23
2	A	301	HEM	C2B-C1B-NB	-2.26	107.24	109.84
2	A	301	HEM	CAA-CBA-CGA	-2.15	107.96	113.67
2	A	301	HEM	CHD-C1D-ND	2.13	126.71	124.42
2	A	301	HEM	CHB-C4A-C3A	-2.07	121.41	127.43
2	A	301	HEM	CHC-C4B-NB	2.03	126.61	124.42
2	A	301	HEM	O1A-CGA-CBA	-2.02	116.68	123.09

There are no chirality outliers.

All (3) torsion outliers are listed below:

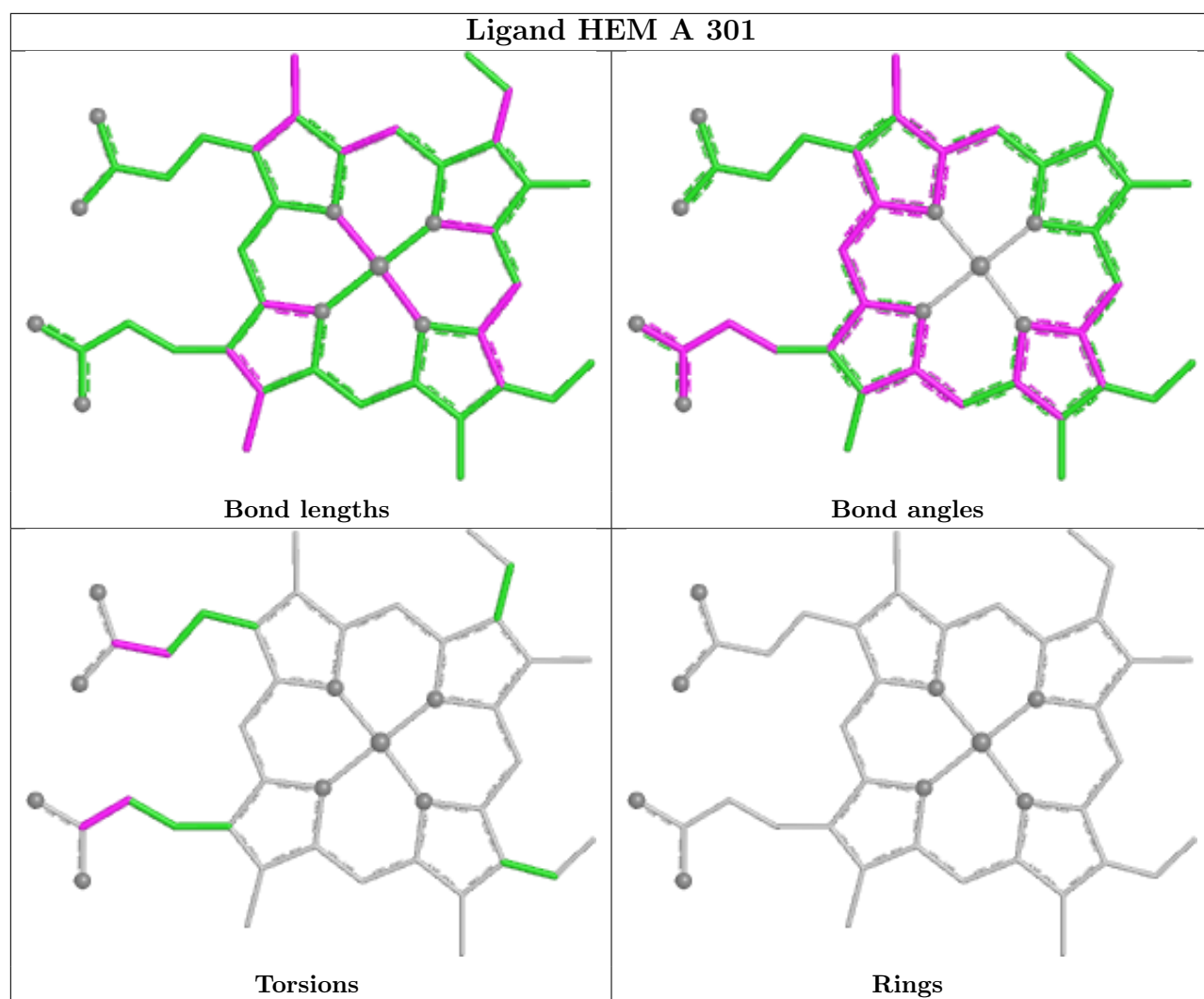
Mol	Chain	Res	Type	Atoms
2	A	301	HEM	CAA-CBA-CGA-O1A
2	A	301	HEM	CAA-CBA-CGA-O2A
2	A	301	HEM	CAD-CBD-CGD-O2D

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	303	SO4	2	0
4	A	304	SO4	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 failed to run properly - this section is therefore empty.

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

EDS failed to run properly - this section is therefore empty.

6.3 Carbohydrates

EDS failed to run properly - this section is therefore empty.

6.4 Ligands

EDS failed to run properly - this section is therefore empty.

6.5 Other polymers

EDS failed to run properly - this section is therefore empty.