



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 12:56 AM UTC

PDB ID : 6NSZ / pdb_00006nsz
Title : X-ray reduced Catalase 3 from N.Crassa (0.526 MGy)
Authors : Zarate-Romero, A.; Rudino-Pinera, E.; Stojanoff, V.
Deposited on : 2019-01-27
Resolution : 2.20 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 : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

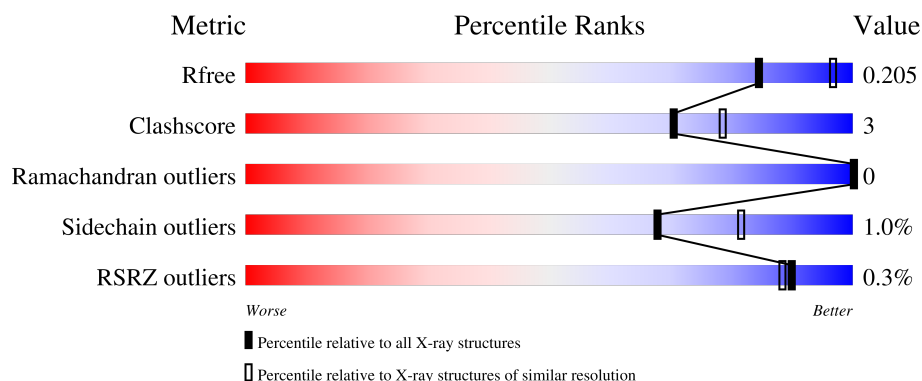
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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))
R_{free}	180053	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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 $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	719	 87% 7% 5%
1	B	719	 86% 8% 6%
1	C	719	 87% 8% 6%
1	D	719	 88% 6% 6%

2 Entry composition [i](#)

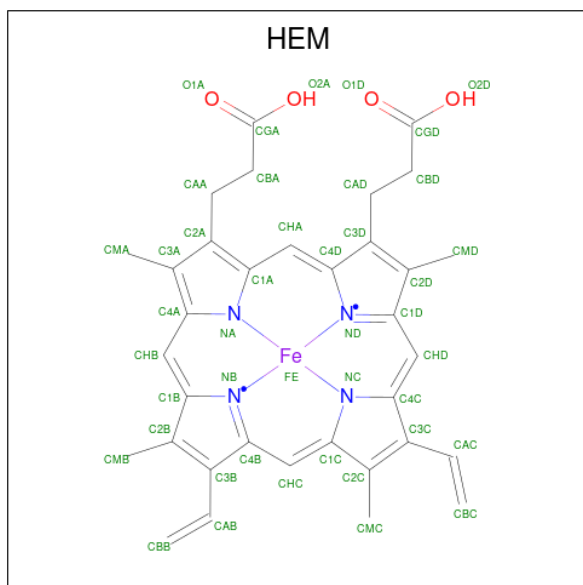
There are 10 unique types of molecules in this entry. The entry contains 23961 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Catalase-3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	682	Total	C	N	O	S	0	9	0
			5420	3426	960	1028	6			
1	B	677	Total	C	N	O	S	0	9	0
			5372	3400	953	1013	6			
1	C	679	Total	C	N	O	S	0	11	0
			5411	3420	962	1023	6			
1	D	678	Total	C	N	O	S	0	9	0
			5385	3408	956	1015	6			

- Molecule 2 is PROTOPORPHYRIN IX CONTAINING FE (CCD ID: HEM) (formula: $C_{34}H_{32}FeN_4O_4$) (labeled as "Ligand of Interest" by depositor).



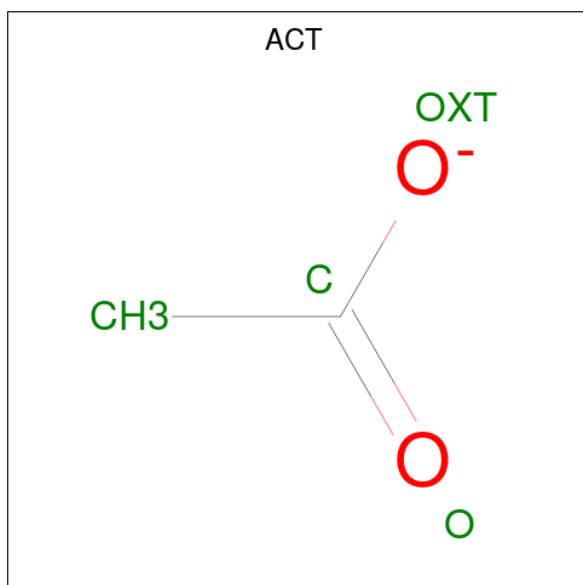
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	Fe	N O	0	0
			43	34	1	4 4		
2	B	1	Total	C	Fe	N O	0	0
			43	34	1	4 4		

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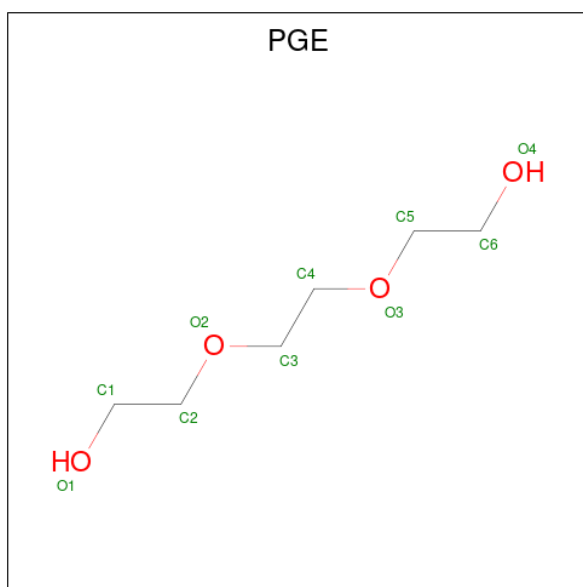
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	C	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
2	D	1	Total 43	C 34	Fe 1	N 4	O 4	0	0

- Molecule 3 is ACETATE ION (CCD ID: ACT) (formula: $C_2H_3O_2$).



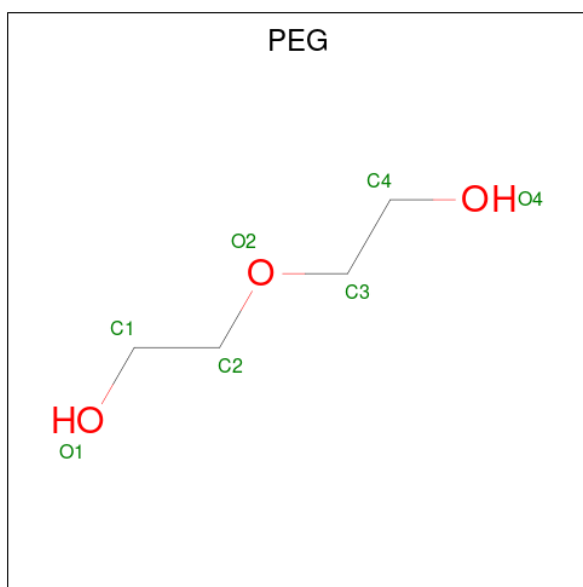
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			4	2	2		
3	A	1	Total	C	O	0	0
			4	2	2		
3	B	1	Total	C	O	0	0
			4	2	2		
3	B	1	Total	C	O	0	0
			4	2	2		
3	C	1	Total	C	O	0	0
			4	2	2		
3	C	1	Total	C	O	0	0
			4	2	2		
3	D	1	Total	C	O	0	0
			4	2	2		
3	D	1	Total	C	O	0	0
			4	2	2		

- Molecule 4 is TRIETHYLENE GLYCOL (CCD ID: PGE) (formula: $C_6H_{14}O_4$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			10	6	4		
4	B	1	Total	C	O	0	0
			10	6	4		
4	D	1	Total	C	O	0	0
			10	6	4		

- Molecule 5 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: $C_4H_{10}O_3$).



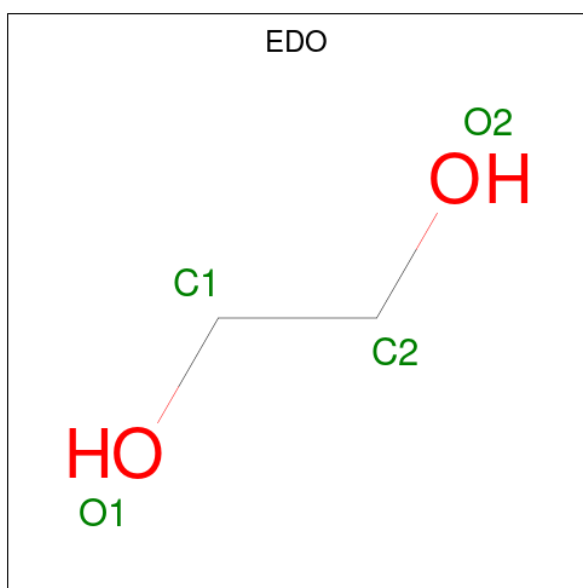
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			7	4	3		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			7	4	3		
5	B	1	Total	C	O	0	0
			7	4	3		
5	C	1	Total	C	O	0	0
			7	4	3		
5	C	1	Total	C	O	0	0
			7	4	3		
5	C	1	Total	C	O	0	0
			7	4	3		

- Molecule 6 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $C_2H_6O_2$).



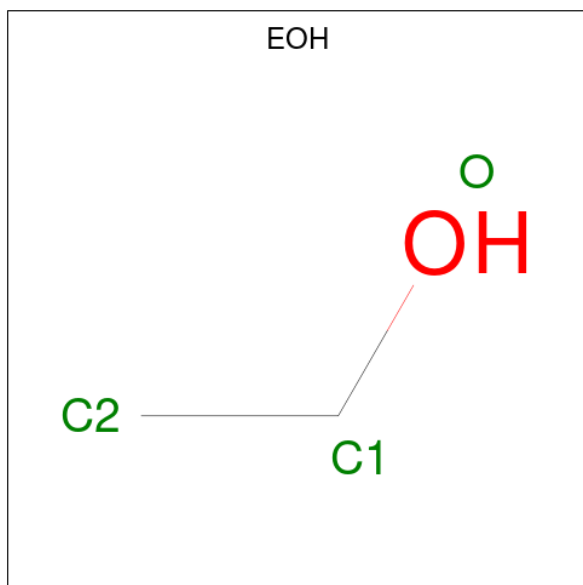
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			4	2	2		
6	B	1	Total	C	O	0	0
			4	2	2		
6	B	1	Total	C	O	0	0
			4	2	2		
6	C	1	Total	C	O	0	0
			4	2	2		
6	C	1	Total	C	O	0	0
			4	2	2		
6	C	1	Total	C	O	0	0
			4	2	2		

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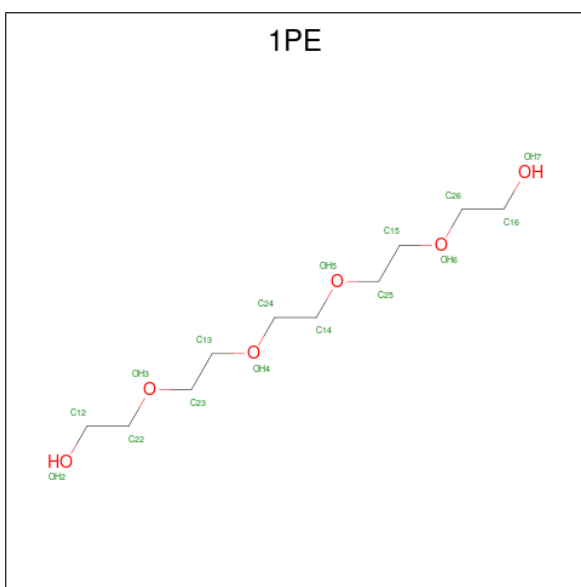
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	D	1	Total	C	O	0	0
			4	2	2		
6	D	1	Total	C	O	0	0
			4	2	2		
6	D	1	Total	C	O	0	0
			4	2	2		

- Molecule 7 is ETHANOL (CCD ID: EOH) (formula: C_2H_6O).



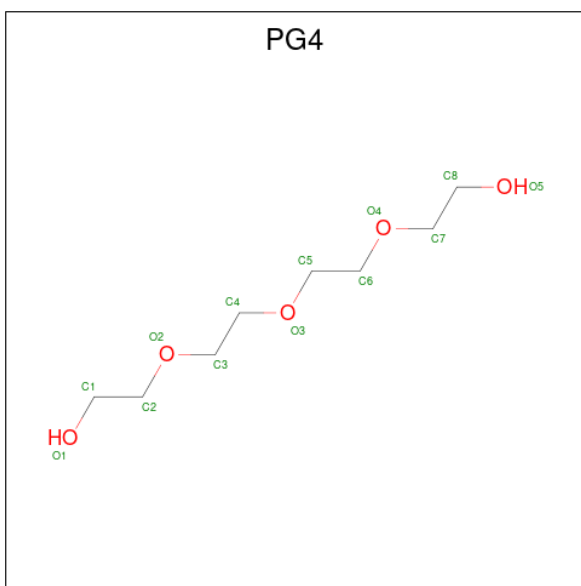
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	C	O	0	0
			3	2	1		
7	B	1	Total	C	O	0	0
			3	2	1		
7	B	1	Total	C	O	0	0
			3	2	1		
7	D	1	Total	C	O	0	0
			3	2	1		
7	D	1	Total	C	O	0	0
			3	2	1		
7	D	1	Total	C	O	0	0
			3	2	1		

- Molecule 8 is PENTAETHYLENE GLYCOL (CCD ID: 1PE) (formula: $C_{10}H_{22}O_6$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	B	1	Total	C	O	0	0
			16	10	6		

- Molecule 9 is TETRAETHYLENE GLYCOL (CCD ID: PG4) (formula: $C_8H_{18}O_5$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
9	C	1	Total	C	O	0	0
			13	8	5		

- Molecule 10 is water.

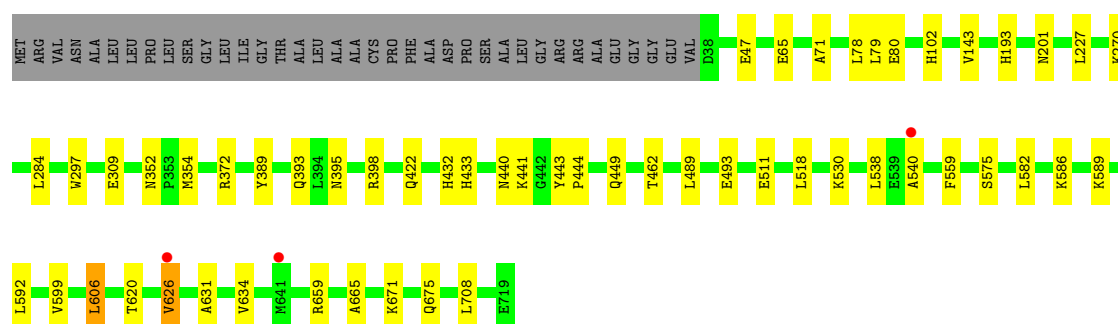
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	A	492	Total 492	O 492	0	0
10	B	494	Total 494	O 494	0	0
10	C	522	Total 522	O 522	0	0
10	D	506	Total 506	O 506	0	0

3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). 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.

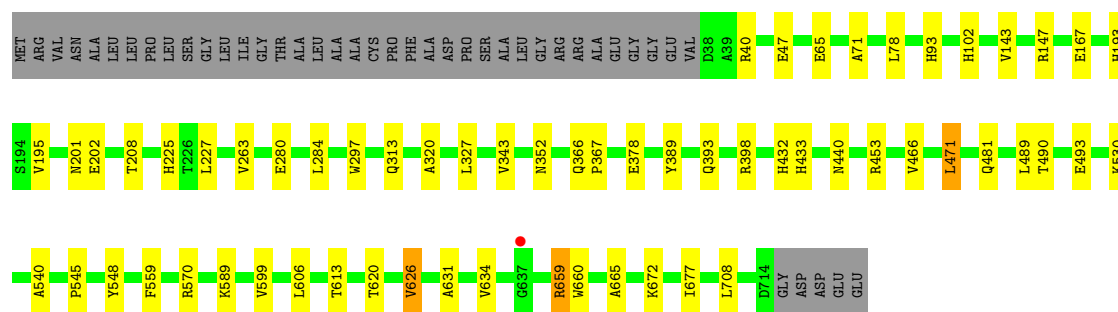
• Molecule 1: Catalase-3

Chain A: 



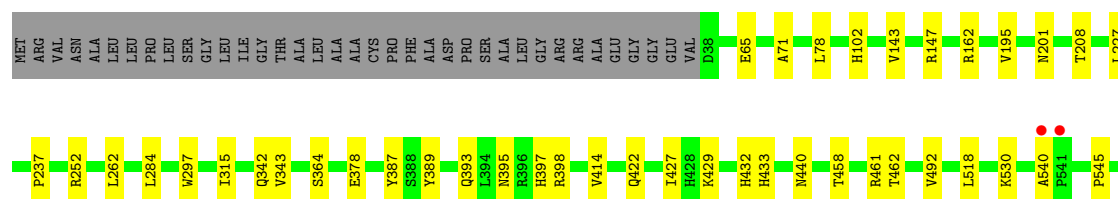
• Molecule 1: Catalase-3

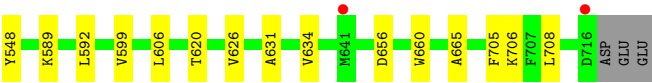
Chain B: 



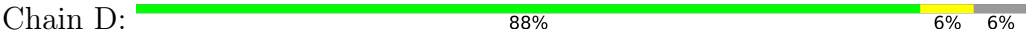
• Molecule 1: Catalase-3

Chain C: 





● Molecule 1: Catalase-3



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	131.10Å 154.50Å 160.90Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.60 – 2.20 29.60 – 2.20	Depositor EDS
% Data completeness (in resolution range)	96.9 (29.60-2.20) 96.8 (29.60-2.20)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.28 (at 2.20Å)	Xtriage
Refinement program	PHENIX 1.12 _2829	Depositor
R, R_{free}	0.157 , 0.204 0.157 , 0.205	Depositor DCC
R_{free} test set	8044 reflections (4.86%)	wwPDB-VP
Wilson B-factor (Å ²)	15.1	Xtriage
Anisotropy	0.025	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 47.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.017 for -h,l,k	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	23961	wwPDB-VP
Average B, all atoms (Å ²)	16.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.76% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: HEM, PGE, PEG, 1PE, ACT, EDO, PG4, EOH

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	0.12	0/5554	0.33	0/7532
1	B	0.12	0/5508	0.34	0/7470
1	C	0.11	0/5544	0.33	0/7517
1	D	0.12	0/5521	0.33	0/7485
All	All	0.12	0/22127	0.33	0/30004

There are no bond length outliers.

There are no bond angle outliers.

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	5420	0	5227	47	0
1	B	5372	0	5198	42	0
1	C	5411	0	5227	39	0
1	D	5385	0	5213	40	0
2	A	43	0	30	1	0
2	B	43	0	30	1	0
2	C	43	0	30	1	0
2	D	43	0	30	2	0
3	A	8	0	6	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	8	0	6	0	0
3	C	8	0	6	2	0
3	D	8	0	6	2	0
4	A	10	0	14	0	0
4	B	10	0	14	0	0
4	D	10	0	14	1	0
5	A	14	0	20	3	0
5	B	7	0	10	1	0
5	C	21	0	30	1	0
6	A	4	0	6	0	0
6	B	8	0	12	0	0
6	C	12	0	18	1	0
6	D	12	0	18	0	0
7	A	3	0	6	0	0
7	B	6	0	12	0	0
7	D	9	0	18	0	0
8	B	16	0	22	1	0
9	C	13	0	18	2	0
10	A	492	0	0	1	0
10	B	494	0	0	3	0
10	C	522	0	0	6	0
10	D	506	0	0	2	0
All	All	23961	0	21241	147	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 3.

All (147) 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:398:ARG:HD2	1:D:398:ARG:CZ	2.00	0.92
1:A:398:ARG:CD	1:D:398:ARG:CZ	2.51	0.88
1:A:398:ARG:NE	1:D:398:ARG:NE	2.21	0.87
1:A:398:ARG:HD2	1:D:398:ARG:NH1	1.90	0.87
1:D:659:ARG:HH12	4:D:804:PGE:H3	1.41	0.86
1:B:398:ARG:NH1	10:B:901:HOH:O	2.17	0.77
1:A:398:ARG:NE	1:D:398:ARG:HE	1.80	0.76
1:C:398:ARG:NH1	10:C:902:HOH:O	2.20	0.73
1:C:530:LYS:HG2	1:C:540:ALA:HB1	1.71	0.73
1:D:398:ARG:HG2	1:D:398:ARG:HH21	1.59	0.68
1:C:342:GLN:HB3	9:C:804:PG4:H62	1.76	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:530:LYS:HG2	1:A:540:ALA:HB1	1.74	0.67
1:C:147[A]:ARG:O	10:C:901:HOH:O	2.13	0.67
1:B:40:ARG:NH2	1:B:47:GLU:OE1	2.27	0.67
1:A:398:ARG:CD	1:D:398:ARG:NE	2.59	0.66
1:D:261:LYS:NZ	10:D:903:HOH:O	2.30	0.65
1:A:80:GLU:OE2	1:C:461[A]:ARG:NH2	2.29	0.65
1:D:398:ARG:HG2	1:D:398:ARG:NH2	2.12	0.65
1:C:71:ALA:HB2	1:C:78:LEU:HD21	1.79	0.64
1:B:570[B]:ARG:NH2	10:B:905:HOH:O	2.31	0.64
1:B:659:ARG:HH12	8:B:804:1PE:H221	1.64	0.62
2:C:801:HEM:HMB2	2:C:801:HEM:HBB2	1.82	0.61
1:A:398:ARG:CD	1:D:398:ARG:NH1	2.64	0.60
9:C:804:PG4:H32	10:C:1291:HOH:O	2.01	0.60
1:B:71:ALA:HB2	1:B:78:LEU:HD21	1.83	0.60
1:B:489:LEU:HB3	1:B:493:GLU:HB3	1.84	0.59
1:A:441:LYS:NZ	5:A:806:PEG:H31	2.17	0.59
2:B:801:HEM:HBB2	2:B:801:HEM:HMB2	1.84	0.59
1:C:195[A]:VAL:HG21	3:C:802:ACT:H1	1.85	0.58
1:C:237:PRO:HB3	3:C:803:ACT:H1	1.86	0.58
1:A:71:ALA:HB2	1:A:78:LEU:HD21	1.87	0.56
1:D:195[A]:VAL:HG21	3:D:802:ACT:H1	1.88	0.56
1:C:147[B]:ARG:NH1	10:C:906:HOH:O	2.29	0.55
1:B:147[A]:ARG:O	10:B:902:HOH:O	2.18	0.54
1:A:354:MET:SD	1:C:429:LYS:HG2	2.48	0.54
1:D:38:ASP:HA	1:D:351:ARG:NH1	2.24	0.53
1:D:237:PRO:HB3	3:D:803:ACT:H1	1.89	0.53
1:A:398:ARG:NE	1:D:398:ARG:CZ	2.69	0.53
1:A:626[B]:VAL:HG22	1:A:665:ALA:HB3	1.90	0.52
1:B:606:LEU:CD1	1:B:613:THR:HG23	2.39	0.52
1:A:270:LYS:HD3	1:A:309:GLU:HG3	1.92	0.52
1:A:589:LYS:HG3	1:A:599:VAL:HB	1.91	0.52
1:D:248:ILE:HD13	2:D:801:HEM:HMB1	1.91	0.52
1:B:320:ALA:HA	1:B:327:LEU:HD12	1.91	0.51
1:A:422:GLN:NE2	10:A:912:HOH:O	2.43	0.51
1:D:492:VAL:HG11	1:D:708:LEU:HB3	1.92	0.51
1:C:458:THR:HB	1:C:461[A]:ARG:HG3	1.93	0.51
1:A:398:ARG:CZ	1:D:398:ARG:HE	2.23	0.51
1:A:398:ARG:CZ	1:D:398:ARG:NE	2.72	0.51
1:C:626:VAL:HG12	1:C:665:ALA:HB3	1.93	0.51
1:B:545:PRO:HA	1:B:548:TYR:CD2	2.46	0.51
1:B:559:PHE:HB3	1:B:708:LEU:HD11	1.92	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:462:THR:HG22	1:D:466:VAL:HG12	1.94	0.50
1:A:441:LYS:HZ2	5:A:806:PEG:H31	1.76	0.49
1:A:441:LYS:HE2	5:A:806:PEG:H22	1.94	0.49
1:C:705:PHE:CD2	1:C:706:LYS:HG2	2.48	0.48
1:D:398:ARG:NH1	10:D:912:HOH:O	2.40	0.48
1:A:671:LYS:O	1:A:675:GLN:HG3	2.13	0.48
1:C:398:ARG:NH2	10:C:907:HOH:O	2.33	0.48
2:D:801:HEM:HMB1	2:D:801:HEM:HBB2	1.96	0.48
1:B:432:HIS:HB2	1:B:440:ASN:HB3	1.94	0.48
1:B:93:HIS:ND1	1:C:395:ASN:HB2	2.28	0.48
1:B:102:HIS:CE1	1:B:143:VAL:HG22	2.49	0.47
1:A:389:TYR:O	1:A:393:GLN:HG2	2.15	0.47
1:C:389:TYR:O	1:C:393:GLN:HG2	2.14	0.47
1:C:432:HIS:HB2	1:C:440:ASN:HB3	1.96	0.47
1:D:389:TYR:O	1:D:393:GLN:HG2	2.14	0.47
1:A:79:LEU:HD11	1:C:387:TYR:HB2	1.96	0.47
1:B:589:LYS:HG3	1:B:599:VAL:HB	1.96	0.47
1:A:582:LEU:O	1:A:586:LYS:HG3	2.15	0.47
1:D:262:LEU:HB2	1:D:315:ILE:HG12	1.97	0.46
1:A:462:THR:HG22	1:B:466:VAL:HG22	1.96	0.46
1:A:422:GLN:NE2	1:C:397:HIS:HA	2.30	0.46
1:D:656:ASP:HB3	1:D:660:TRP:CZ3	2.51	0.46
1:B:284:LEU:HD13	1:B:620:THR:HB	1.96	0.46
1:C:656:ASP:HB3	1:C:660:TRP:CZ3	2.51	0.46
1:C:589:LYS:HG3	1:C:599:VAL:HB	1.98	0.46
1:C:545:PRO:HA	1:C:548:TYR:CD2	2.51	0.46
1:B:389:TYR:O	1:B:393:GLN:HG2	2.16	0.46
1:B:626:VAL:HB	1:B:665:ALA:HB3	1.97	0.46
1:D:71:ALA:HB2	1:D:78:LEU:HD21	1.98	0.46
1:C:252:ARG:NH1	1:C:461[A]:ARG:HD3	2.31	0.46
1:C:414:VAL:HB	6:C:808:EDO:H22	1.98	0.46
1:C:162:ARG:NH2	1:C:364:SER:O	2.45	0.45
1:C:378:GLU:N	1:C:378:GLU:OE1	2.49	0.45
1:B:195[B]:VAL:HG13	1:B:208:THR:HG22	1.98	0.45
1:C:284:LEU:HD13	1:C:620:THR:HB	1.99	0.45
1:A:102:HIS:CE1	1:A:143:VAL:HG22	2.51	0.45
1:C:102:HIS:CE1	1:C:143:VAL:HG22	2.51	0.45
1:A:78:LEU:HD11	1:D:467:LEU:HD22	1.99	0.45
1:B:167:GLU:OE2	1:B:453:ARG:NE	2.44	0.45
1:B:297:TRP:CZ3	1:B:352:ASN:HB3	2.52	0.45
1:B:378:GLU:N	1:B:378:GLU:OE1	2.50	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:378:GLU:N	1:D:378:GLU:OE1	2.50	0.45
1:B:606:LEU:HD13	1:B:613:THR:HG23	1.98	0.45
1:B:367:PRO:HD3	1:B:389:TYR:CD2	2.52	0.45
1:C:492:VAL:HG11	1:C:708:LEU:HB3	1.97	0.45
1:A:432:HIS:HB2	1:A:440:ASN:HB3	1.98	0.44
1:B:677:ILE:O	5:B:806:PEG:H32	2.17	0.44
1:D:545:PRO:HA	1:D:548:TYR:CD2	2.53	0.44
1:A:193:HIS:ND1	1:B:433:HIS:CD2	2.86	0.44
1:A:511:GLU:OE2	1:A:538:LEU:HD21	2.17	0.44
1:A:631:ALA:O	1:A:634:VAL:HG22	2.17	0.44
1:B:398:ARG:HD3	1:D:422:GLN:NE2	2.32	0.44
1:A:575:SER:HA	1:A:631:ALA:HB2	1.99	0.43
1:A:433:HIS:CE1	1:B:193:HIS:HB3	2.54	0.43
1:C:631:ALA:O	1:C:634:VAL:HG22	2.18	0.43
1:D:280:GLU:HB2	1:D:660:TRP:CD2	2.52	0.43
1:D:631:ALA:O	1:D:634:VAL:HG22	2.17	0.43
1:B:201:ASN:HB3	1:C:297:TRP:CE3	2.53	0.43
1:C:518[A]:LEU:HD23	1:C:518[A]:LEU:HA	1.92	0.43
1:D:681:GLU:H	1:D:681:GLU:HG3	1.67	0.43
1:A:395:ASN:OD1	1:A:398:ARG:NH1	2.51	0.43
1:A:398:ARG:HG3	1:C:422:GLN:NE2	2.33	0.43
1:A:559:PHE:HB3	1:A:708:LEU:HD11	2.00	0.43
1:A:489:LEU:HB3	1:A:493:GLU:HB3	2.00	0.43
1:A:606:LEU:HD12	1:A:606:LEU:HA	1.89	0.43
1:B:280:GLU:HB2	1:B:660:TRP:CD2	2.54	0.43
1:B:297:TRP:CE3	1:C:201:ASN:HB3	2.54	0.43
1:B:490:THR:OG1	1:B:493:GLU:HB2	2.19	0.43
1:B:202:GLU:OE2	1:C:147[A]:ARG:NH2	2.50	0.42
1:B:631:ALA:O	1:B:634:VAL:HG22	2.18	0.42
1:C:262:LEU:HB2	1:C:315:ILE:HB	2.01	0.42
1:A:297:TRP:CZ3	1:A:352:ASN:HB3	2.55	0.42
2:A:801:HEM:HMB1	2:A:801:HEM:HBB2	2.00	0.42
1:A:443:TYR:HA	1:A:444:PRO:C	2.45	0.42
1:A:47:GLU:HA	1:C:427:ILE:O	2.20	0.41
1:A:201:ASN:HB3	1:D:297:TRP:CE3	2.55	0.41
1:A:284:LEU:HD13	1:A:620:THR:HB	2.02	0.41
1:B:606:LEU:HD13	1:B:613:THR:CG2	2.50	0.41
1:D:102:HIS:CE1	1:D:143:VAL:HG22	2.55	0.41
1:C:195[B]:VAL:HG13	1:C:208:THR:HG22	2.01	0.41
1:D:575:SER:HA	1:D:631:ALA:HB2	2.02	0.41
1:A:518:LEU:HD23	1:A:518:LEU:HA	1.89	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:366:GLN:HG2	1:B:389:TYR:HB3	2.01	0.41
1:A:372:ARG:HH12	1:A:449:GLN:HA	1.84	0.41
1:D:284:LEU:HD13	1:D:620:THR:HB	2.03	0.41
1:B:471:LEU:HD12	1:B:471:LEU:HA	1.86	0.41
1:B:530:LYS:HA	1:B:540:ALA:HB2	2.03	0.41
1:D:385:ARG:O	1:D:389:TYR:HD1	2.04	0.41
1:D:682:LYS:HE2	1:D:682:LYS:HB3	1.92	0.41
1:B:225:HIS:CE1	1:B:481:GLN:HB3	2.56	0.41
1:B:672:LYS:HA	1:B:672:LYS:HD2	1.98	0.41
1:D:589:LYS:HG3	1:D:599:VAL:HB	2.03	0.41
1:B:263:VAL:HA	1:B:313:GLN:O	2.21	0.40
5:C:806:PEG:H22	10:C:1137:HOH:O	2.22	0.40
1:D:471:LEU:HD12	1:D:471:LEU:HA	1.92	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	689/719 (96%)	667 (97%)	22 (3%)	0	100	100
1	B	683/719 (95%)	663 (97%)	20 (3%)	0	100	100
1	C	688/719 (96%)	666 (97%)	22 (3%)	0	100	100
1	D	685/719 (95%)	666 (97%)	19 (3%)	0	100	100
All	All	2745/2876 (95%)	2662 (97%)	83 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	568/585 (97%)	561 (99%)	7 (1%)	63	78
1	B	563/585 (96%)	557 (99%)	6 (1%)	65	79
1	C	567/585 (97%)	561 (99%)	6 (1%)	65	79
1	D	564/585 (96%)	560 (99%)	4 (1%)	76	87
All	All	2262/2340 (97%)	2239 (99%)	23 (1%)	68	81

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	65	GLU
1	A	227	LEU
1	A	592	LEU
1	A	606	LEU
1	A	626[A]	VAL
1	A	626[B]	VAL
1	A	659	ARG
1	B	65	GLU
1	B	227	LEU
1	B	343	VAL
1	B	471	LEU
1	B	626	VAL
1	B	659	ARG
1	C	65	GLU
1	C	227	LEU
1	C	343	VAL
1	C	433	HIS
1	C	592	LEU
1	C	606	LEU
1	D	65	GLU
1	D	227	LEU
1	D	315	ILE
1	D	709	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	133	GLN
1	A	183	GLN
1	A	220	GLN
1	A	422	GLN
1	A	520	GLN
1	A	522	ASN
1	A	526	ASN
1	B	133	GLN
1	B	512	GLN
1	B	520	GLN
1	B	652	GLN
1	C	220	GLN
1	C	520	GLN
1	C	550	ASN
1	C	591	GLN
1	C	652	GLN
1	D	183	GLN
1	D	220	GLN
1	D	526	ASN
1	D	612	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](#)

38 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
6	EDO	D	806	-	3,3,3	0.43	0	2,2,2	0.41	0
5	PEG	B	806	-	6,6,6	0.51	0	5,5,5	0.40	0
3	ACT	B	803	-	3,3,3	1.64	1 (33%)	3,3,3	1.15	0
4	PGE	D	804	-	9,9,9	0.36	0	8,8,8	0.22	0
4	PGE	A	804	-	9,9,9	0.31	0	8,8,8	0.33	0
2	HEM	D	801	10,1	50,50,50	1.42	8 (16%)	67,82,82	1.17	6 (8%)
7	EOH	A	808	-	2,2,2	0.47	0	1,1,1	0.14	0
8	1PE	B	804	-	15,15,15	0.54	0	14,14,14	0.40	0
3	ACT	C	803	-	3,3,3	1.61	1 (33%)	3,3,3	1.55	0
6	EDO	B	807	-	3,3,3	0.41	0	2,2,2	0.31	0
3	ACT	A	802	-	3,3,3	1.44	0	3,3,3	1.42	0
3	ACT	A	803	-	3,3,3	1.73	1 (33%)	3,3,3	1.24	0
9	PG4	C	804	-	12,12,12	0.52	0	11,11,11	0.44	0
5	PEG	A	805	-	6,6,6	0.48	0	5,5,5	0.54	0
6	EDO	C	808	-	3,3,3	0.46	0	2,2,2	0.29	0
7	EOH	B	809	-	2,2,2	0.44	0	1,1,1	0.18	0
6	EDO	A	807	-	3,3,3	0.44	0	2,2,2	0.32	0
5	PEG	C	806	-	6,6,6	0.49	0	5,5,5	0.44	0
6	EDO	D	805	-	3,3,3	0.49	0	2,2,2	0.34	0
5	PEG	C	805	-	6,6,6	0.50	0	5,5,5	0.42	0
6	EDO	C	809	-	3,3,3	0.43	0	2,2,2	0.42	0
2	HEM	B	801	1	50,50,50	1.44	9 (18%)	67,82,82	1.13	5 (7%)
5	PEG	A	806	-	6,6,6	0.45	0	5,5,5	0.53	0
7	EOH	D	810	-	2,2,2	0.43	0	1,1,1	0.16	0
7	EOH	B	810	-	2,2,2	0.47	0	1,1,1	0.20	0
7	EOH	D	808	-	2,2,2	0.46	0	1,1,1	0.15	0
3	ACT	D	803	-	3,3,3	1.70	1 (33%)	3,3,3	1.28	0
3	ACT	B	802	-	3,3,3	1.48	1 (33%)	3,3,3	1.31	0
3	ACT	C	802	-	3,3,3	1.45	1 (33%)	3,3,3	1.37	0
3	ACT	D	802	-	3,3,3	1.23	0	3,3,3	1.51	0
6	EDO	B	808	-	3,3,3	0.45	0	2,2,2	0.39	0
2	HEM	C	801	10,1	50,50,50	1.43	8 (16%)	67,82,82	1.16	6 (8%)
5	PEG	C	807	-	6,6,6	0.51	0	5,5,5	0.48	0
6	EDO	D	807	-	3,3,3	0.46	0	2,2,2	0.32	0
2	HEM	A	801	10,1	50,50,50	1.42	9 (18%)	67,82,82	1.14	5 (7%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	EDO	C	810	-	3,3,3	0.40	0	2,2,2	0.47	0
7	EOH	D	809	-	2,2,2	0.47	0	1,1,1	0.14	0
4	PGE	B	805	-	9,9,9	0.33	0	8,8,8	0.31	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
6	EDO	D	806	-	-	0/1/1/1	-
5	PEG	B	806	-	-	2/4/4/4	-
4	PGE	D	804	-	-	4/7/7/7	-
4	PGE	A	804	-	-	3/7/7/7	-
2	HEM	D	801	10,1	-	2/14/54/54	-
8	1PE	B	804	-	-	8/13/13/13	-
6	EDO	B	807	-	-	0/1/1/1	-
9	PG4	C	804	-	-	7/10/10/10	-
5	PEG	A	805	-	-	2/4/4/4	-
6	EDO	C	808	-	-	0/1/1/1	-
6	EDO	A	807	-	-	0/1/1/1	-
5	PEG	C	806	-	-	0/4/4/4	-
6	EDO	D	805	-	-	1/1/1/1	-
5	PEG	C	805	-	-	2/4/4/4	-
6	EDO	C	809	-	-	1/1/1/1	-
2	HEM	B	801	1	-	4/14/54/54	-
5	PEG	A	806	-	-	1/4/4/4	-
6	EDO	B	808	-	-	0/1/1/1	-
2	HEM	C	801	10,1	-	2/14/54/54	-
5	PEG	C	807	-	-	1/4/4/4	-
6	EDO	D	807	-	-	0/1/1/1	-
2	HEM	A	801	10,1	-	2/14/54/54	-
6	EDO	C	810	-	-	1/1/1/1	-
4	PGE	B	805	-	-	5/7/7/7	-

All (40) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	801	HEM	FE-NB	3.99	2.07	1.94

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	801	HEM	FE-NA	3.86	2.07	1.95
2	D	801	HEM	FE-ND	3.85	2.06	1.94
2	B	801	HEM	FE-ND	3.80	2.06	1.94
2	A	801	HEM	FE-ND	3.57	2.05	1.94
2	A	801	HEM	FE-NB	3.51	2.05	1.94
2	B	801	HEM	FE-NC	3.50	2.06	1.95
2	C	801	HEM	FE-ND	3.20	2.04	1.94
2	B	801	HEM	FE-NB	3.12	2.04	1.94
2	D	801	HEM	CAB-C3B	3.09	1.55	1.47
2	B	801	HEM	CAB-C3B	3.08	1.55	1.47
2	C	801	HEM	CAB-C3B	3.06	1.55	1.47
2	D	801	HEM	CAC-C3C	3.04	1.55	1.47
2	A	801	HEM	CAB-C3B	3.02	1.55	1.47
2	C	801	HEM	FE-NC	3.00	2.05	1.95
2	C	801	HEM	CAC-C3C	3.00	1.55	1.47
2	B	801	HEM	CAC-C3C	3.00	1.55	1.47
2	A	801	HEM	CAC-C3C	3.00	1.55	1.47
2	A	801	HEM	FE-NA	2.93	2.04	1.95
2	D	801	HEM	FE-NB	2.89	2.03	1.94
2	C	801	HEM	FE-NA	2.89	2.04	1.95
2	A	801	HEM	FE-NC	2.71	2.04	1.95
2	B	801	HEM	FE-NA	2.61	2.03	1.95
3	A	803	ACT	CH3-C	2.56	1.59	1.49
3	D	803	ACT	CH3-C	2.52	1.59	1.49
3	B	803	ACT	CH3-C	2.48	1.58	1.49
3	C	803	ACT	CH3-C	2.44	1.58	1.49
2	A	801	HEM	CMB-C2B	2.20	1.55	1.50
2	C	801	HEM	CMB-C2B	2.15	1.55	1.50
2	B	801	HEM	CMB-C2B	2.15	1.55	1.50
3	C	802	ACT	CH3-C	2.12	1.57	1.49
2	B	801	HEM	CMC-C2C	2.11	1.55	1.50
2	D	801	HEM	CMC-C2C	2.10	1.55	1.50
2	D	801	HEM	FE-NC	2.08	2.02	1.95
2	D	801	HEM	CMB-C2B	2.07	1.55	1.50
3	B	802	ACT	CH3-C	2.06	1.57	1.49
2	A	801	HEM	CMD-C2D	2.05	1.55	1.50
2	A	801	HEM	CMC-C2C	2.03	1.54	1.50
2	C	801	HEM	CMC-C2C	2.03	1.54	1.50
2	B	801	HEM	CMA-C3A	2.00	1.54	1.50

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	801	HEM	CAA-CBA-CGA	-2.72	106.46	113.67
2	C	801	HEM	CAA-CBA-CGA	-2.67	106.58	113.67
2	A	801	HEM	CAA-CBA-CGA	-2.65	106.64	113.67
2	D	801	HEM	CAA-CBA-CGA	-2.64	106.65	113.67
2	C	801	HEM	C4D-ND-C1D	2.60	108.28	105.21
2	D	801	HEM	C4D-ND-C1D	2.57	108.25	105.21
2	B	801	HEM	C4D-ND-C1D	2.56	108.23	105.21
2	D	801	HEM	C1B-NB-C4B	2.53	108.20	105.21
2	A	801	HEM	C4D-ND-C1D	2.48	108.15	105.21
2	C	801	HEM	C2A-C1A-NA	-2.44	107.45	110.15
2	C	801	HEM	C1B-NB-C4B	2.44	108.09	105.21
2	B	801	HEM	C1B-NB-C4B	2.40	108.05	105.21
2	A	801	HEM	C1B-NB-C4B	2.37	108.02	105.21
2	A	801	HEM	C2A-C1A-NA	-2.31	107.59	110.15
2	C	801	HEM	C3D-C4D-ND	-2.29	107.66	110.17
2	D	801	HEM	C2A-C1A-NA	-2.29	107.61	110.15
2	B	801	HEM	C2A-C1A-NA	-2.26	107.65	110.15
2	D	801	HEM	C3D-C4D-ND	-2.23	107.73	110.17
2	B	801	HEM	C3D-C4D-ND	-2.22	107.74	110.17
2	A	801	HEM	C3D-C4D-ND	-2.22	107.74	110.17
2	D	801	HEM	C3B-C2B-C1B	2.21	108.07	106.41
2	C	801	HEM	C3B-C2B-C1B	2.16	108.03	106.41

There are no chirality outliers.

All (48) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	804	PGE	O2-C3-C4-O3
8	B	804	1PE	OH5-C14-C24-OH4
8	B	804	1PE	OH6-C15-C25-OH5
9	C	804	PG4	O2-C3-C4-O3
5	C	807	PEG	O1-C1-C2-O2
8	B	804	1PE	OH4-C13-C23-OH3
4	A	804	PGE	O1-C1-C2-O2
4	D	804	PGE	O3-C5-C6-O4
8	B	804	1PE	OH7-C16-C26-OH6
9	C	804	PG4	C5-C6-O4-C7
8	B	804	1PE	C23-C13-OH4-C24
4	D	804	PGE	O1-C1-C2-O2
5	A	806	PEG	O2-C3-C4-O4
8	B	804	1PE	OH2-C12-C22-OH3
4	B	805	PGE	O2-C3-C4-O3
9	C	804	PG4	O3-C5-C6-O4

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Mol	Chain	Res	Type	Atoms
5	C	805	PEG	C4-C3-O2-C2
4	B	805	PGE	C1-C2-O2-C3
5	A	805	PEG	C4-C3-O2-C2
4	B	805	PGE	C4-C3-O2-C2
9	C	804	PG4	C8-C7-O4-C6
4	B	805	PGE	O3-C5-C6-O4
5	B	806	PEG	C1-C2-O2-C3
6	C	810	EDO	O1-C1-C2-O2
6	D	805	EDO	O1-C1-C2-O2
9	C	804	PG4	C6-C5-O3-C4
5	C	805	PEG	C1-C2-O2-C3
5	A	805	PEG	C1-C2-O2-C3
4	B	805	PGE	C6-C5-O3-C4
9	C	804	PG4	C4-C3-O2-C2
9	C	804	PG4	C1-C2-O2-C3
2	A	801	HEM	CAA-CBA-CGA-O1A
2	B	801	HEM	CAA-CBA-CGA-O1A
5	B	806	PEG	C4-C3-O2-C2
2	A	801	HEM	CAA-CBA-CGA-O2A
4	D	804	PGE	O2-C3-C4-O3
2	B	801	HEM	CAA-CBA-CGA-O2A
8	B	804	1PE	C25-C15-OH6-C26
2	C	801	HEM	CAA-CBA-CGA-O2A
4	A	804	PGE	O3-C5-C6-O4
2	D	801	HEM	CAA-CBA-CGA-O1A
2	D	801	HEM	CAA-CBA-CGA-O2A
2	C	801	HEM	CAA-CBA-CGA-O1A
6	C	809	EDO	O1-C1-C2-O2
2	B	801	HEM	CAD-CBD-CGD-O2D
4	D	804	PGE	C6-C5-O3-C4
8	B	804	1PE	C13-C23-OH3-C22
2	B	801	HEM	CAD-CBD-CGD-O1D

There are no ring outliers.

15 monomers are involved in 19 short contacts:

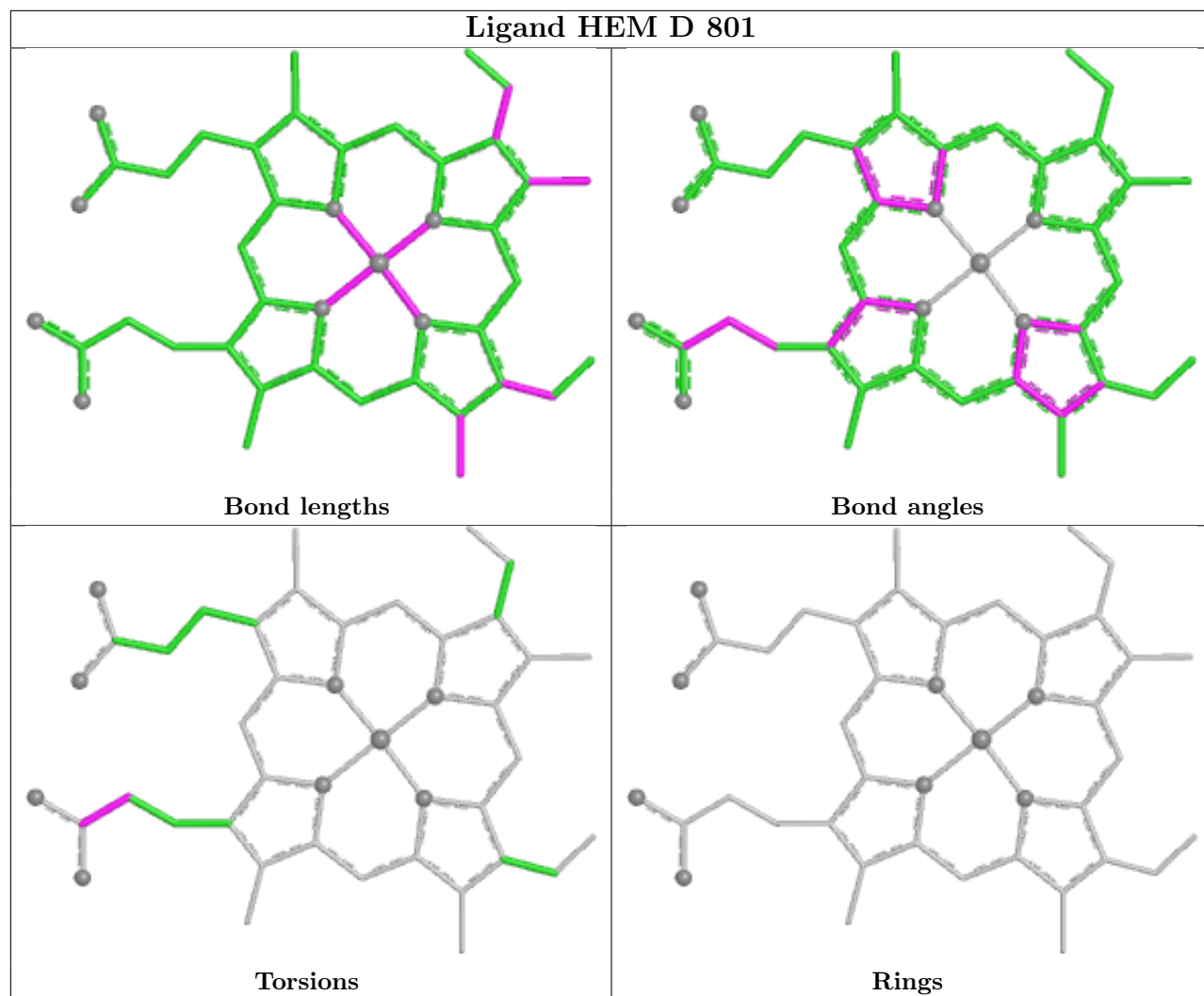
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	806	PEG	1	0
4	D	804	PGE	1	0
2	D	801	HEM	2	0
8	B	804	1PE	1	0
3	C	803	ACT	1	0

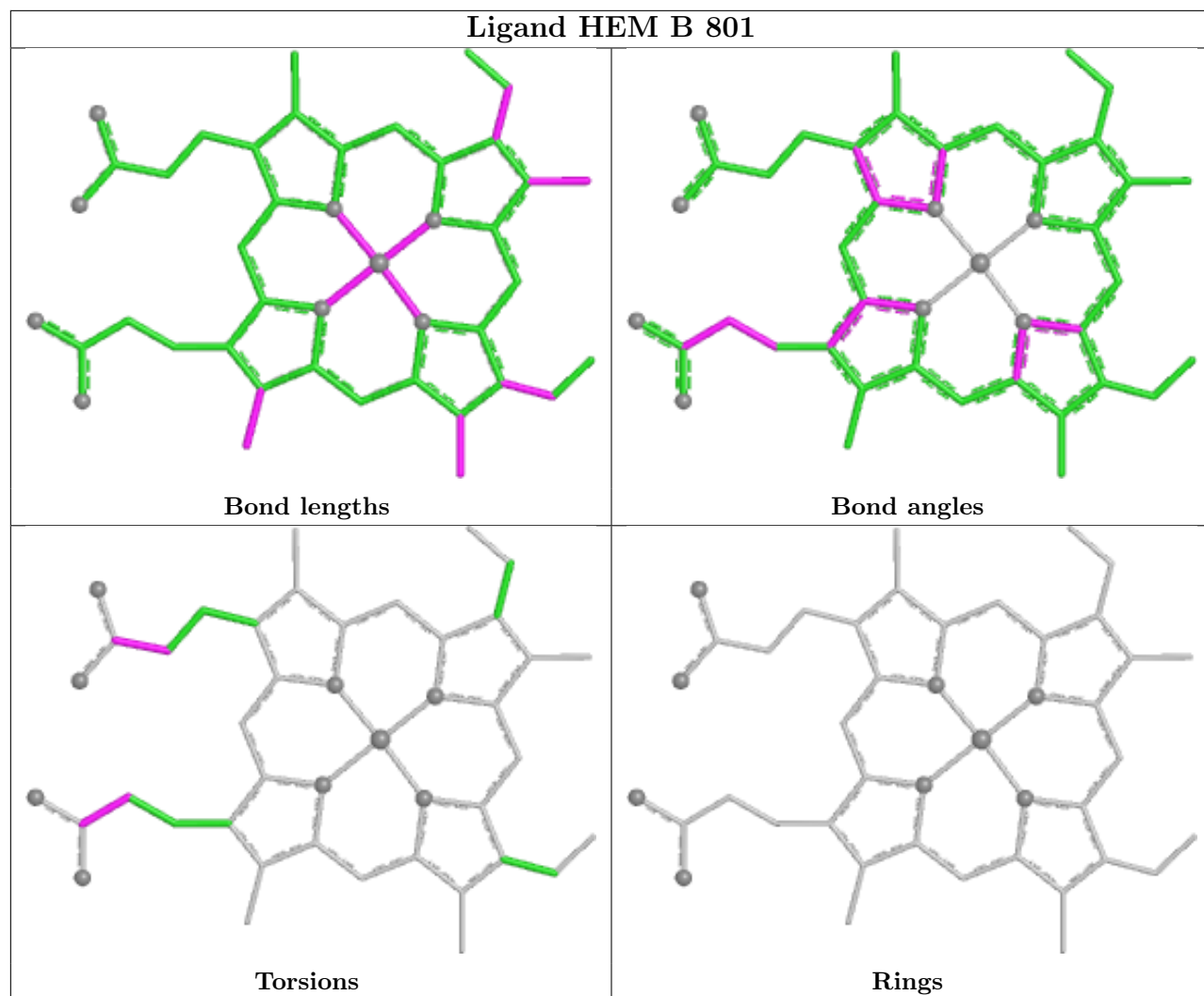
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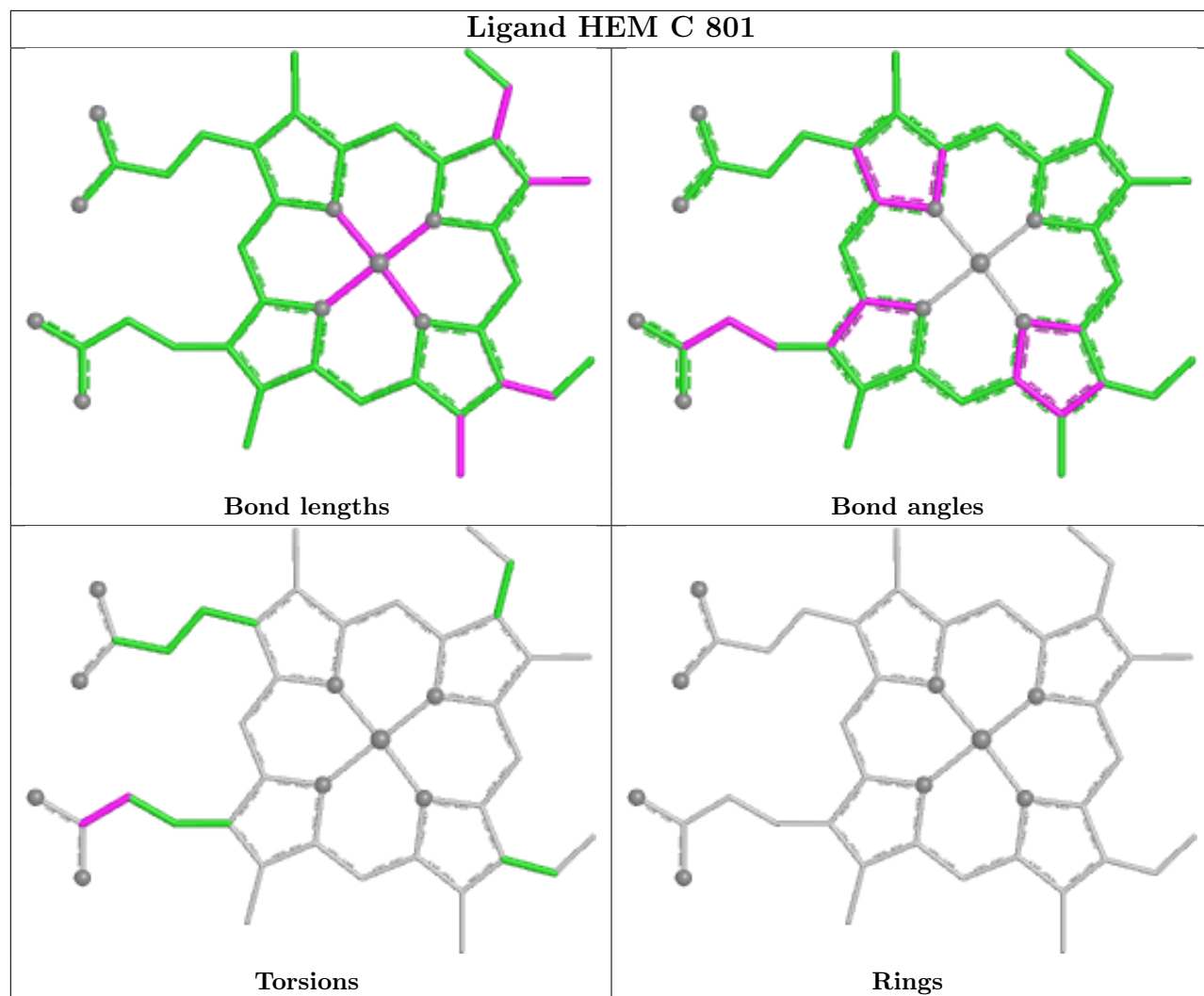
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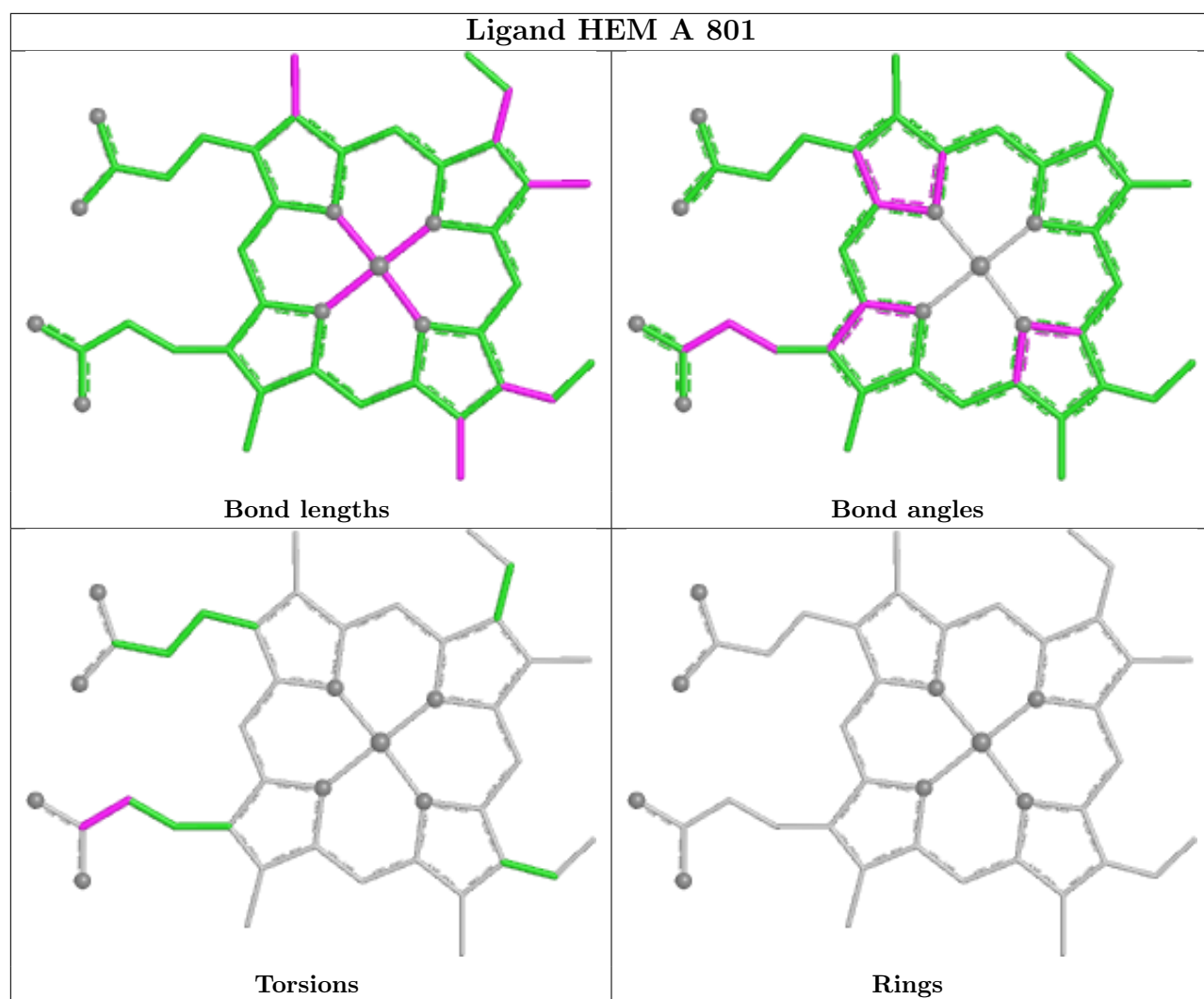
Mol	Chain	Res	Type	Clashes	Symm-Clashes
9	C	804	PG4	2	0
6	C	808	EDO	1	0
5	C	806	PEG	1	0
2	B	801	HEM	1	0
5	A	806	PEG	3	0
3	D	803	ACT	1	0
3	C	802	ACT	1	0
3	D	802	ACT	1	0
2	C	801	HEM	1	0
2	A	801	HEM	1	0

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	682/719 (94%)	-0.48	3 (0%) 88 87	4, 15, 36, 65	9 (1%)
1	B	677/719 (94%)	-0.56	1 (0%) 92 90	5, 13, 31, 55	8 (1%)
1	C	679/719 (94%)	-0.55	4 (0%) 85 83	4, 13, 32, 64	11 (1%)
1	D	678/719 (94%)	-0.51	1 (0%) 92 90	4, 15, 34, 66	9 (1%)
All	All	2716/2876 (94%)	-0.52	9 (0%) 90 88	4, 14, 33, 66	37 (1%)

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	716	ASP	4.4
1	B	637	GLY	3.1
1	C	540	ALA	2.8
1	A	641	MET	2.4
1	D	680	GLU	2.4
1	A	626[A]	VAL	2.4
1	C	641	MET	2.1
1	C	541	PRO	2.1
1	A	540	ALA	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	ACT	D	803	4/4	0.73	0.17	20,21,22,29	0
7	EOH	D	810	3/3	0.73	0.15	22,22,24,27	0
6	EDO	C	809	4/4	0.74	0.14	29,29,30,33	0
7	EOH	B	809	3/3	0.78	0.14	23,23,26,29	0
3	ACT	A	803	4/4	0.79	0.14	19,19,19,23	0
3	ACT	C	803	4/4	0.79	0.15	23,24,25,35	0
3	ACT	B	803	4/4	0.80	0.13	17,17,18,23	0
8	1PE	B	804	16/16	0.80	0.14	30,31,33,34	0
6	EDO	D	806	4/4	0.81	0.12	30,31,33,33	0
6	EDO	B	808	4/4	0.81	0.13	32,32,34,34	0
5	PEG	B	806	7/7	0.81	0.22	24,26,27,28	0
6	EDO	D	805	4/4	0.81	0.15	30,32,32,33	0
7	EOH	D	808	3/3	0.82	0.16	29,29,32,36	0
4	PGE	D	804	10/10	0.84	0.15	30,31,32,32	0
5	PEG	A	805	7/7	0.84	0.12	26,27,30,30	0
7	EOH	B	810	3/3	0.84	0.14	23,23,23,26	0
6	EDO	C	810	4/4	0.85	0.12	27,27,29,29	0
7	EOH	D	809	3/3	0.85	0.10	24,24,25,27	0
4	PGE	A	804	10/10	0.86	0.14	26,27,30,31	0
5	PEG	C	806	7/7	0.87	0.12	27,27,29,30	0
9	PG4	C	804	13/13	0.87	0.11	27,27,29,29	0
6	EDO	D	807	4/4	0.88	0.12	29,30,30,32	0
5	PEG	A	806	7/7	0.89	0.12	20,22,24,26	0
4	PGE	B	805	10/10	0.90	0.12	24,27,30,31	0
6	EDO	C	808	4/4	0.90	0.12	30,30,31,31	0
5	PEG	C	805	7/7	0.91	0.12	30,31,32,33	0
7	EOH	A	808	3/3	0.91	0.10	23,23,24,26	0
6	EDO	B	807	4/4	0.91	0.09	26,27,28,29	0
5	PEG	C	807	7/7	0.93	0.07	26,27,28,29	0
6	EDO	A	807	4/4	0.94	0.08	27,28,29,30	0
3	ACT	B	802	4/4	0.96	0.08	10,12,12,12	0
3	ACT	A	802	4/4	0.96	0.07	15,16,16,16	0
3	ACT	D	802	4/4	0.97	0.06	13,14,15,16	0
2	HEM	B	801	43/43	0.97	0.06	3,8,11,15	0
2	HEM	A	801	43/43	0.98	0.06	4,8,13,17	0
2	HEM	C	801	43/43	0.98	0.05	3,7,11,12	0
2	HEM	D	801	43/43	0.98	0.05	2,7,12,18	0

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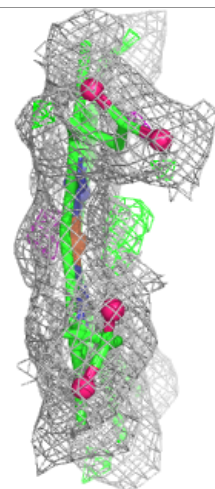
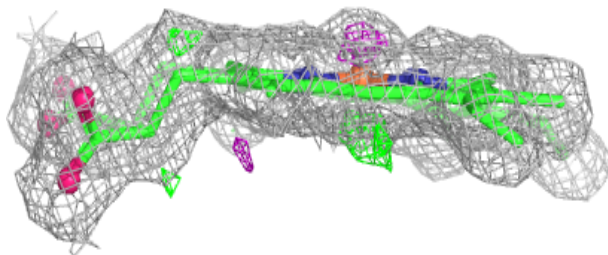
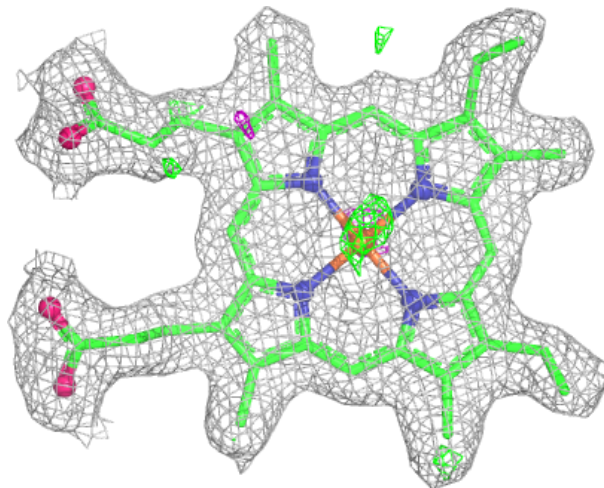
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	ACT	C	802	4/4	0.99	0.04	16,16,17,17	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

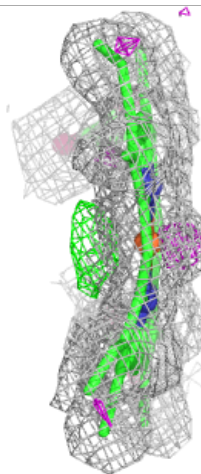
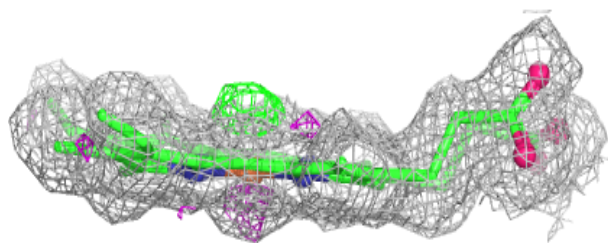
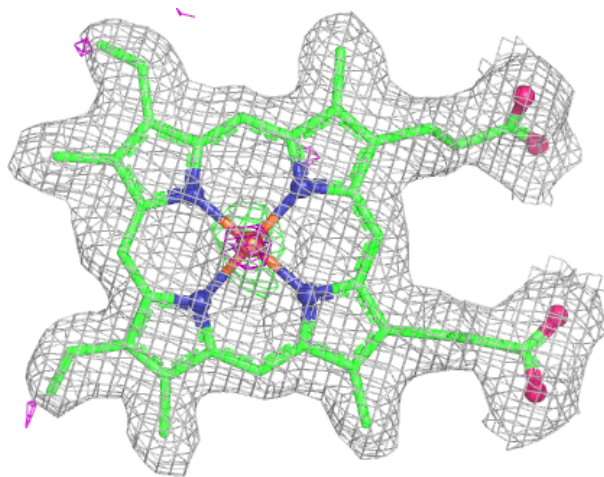
Electron density around HEM B 801:

$2mF_o - DF_c$ (at 0.7 rmsd) in gray
 $mF_o - DF_c$ (at 3 rmsd) in purple (negative)
 and green (positive)



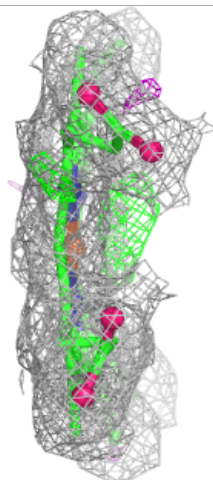
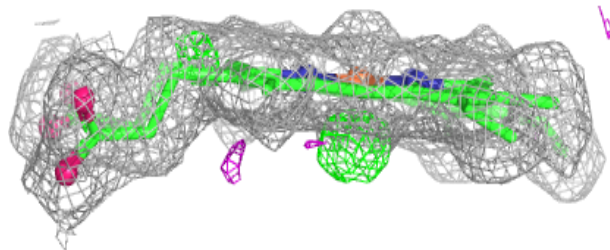
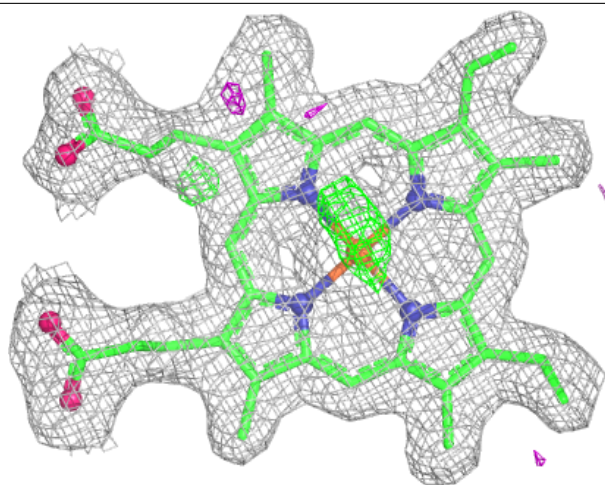
Electron density around HEM A 801:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



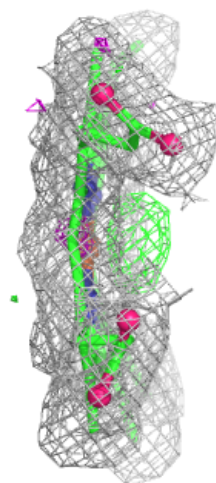
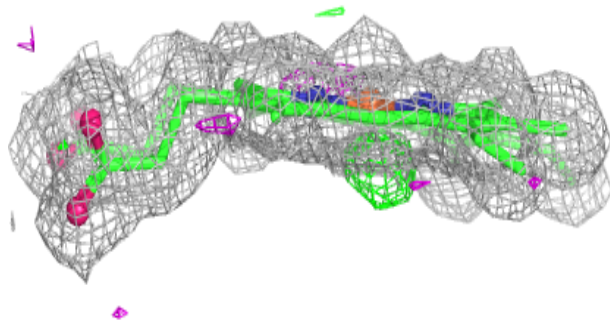
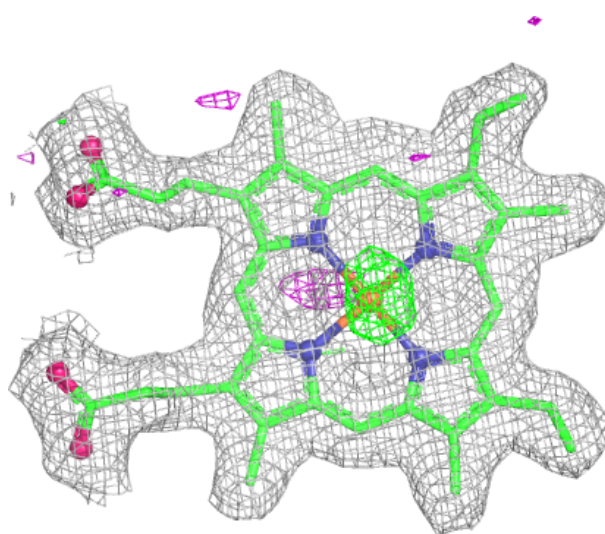
Electron density around HEM C 801:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around HEM D 801:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers ⓘ

There are no such residues in this entry.