



# Full wwPDB X-ray Structure Validation Report ⓘ

Jun 12, 2024 – 09:38 PM EDT

PDB ID : 3ZJ5  
Title : NEUROSPORA CRASSA CATALASE-3 EXPRESSED IN E. COLI, ORTHORHOMBIC FORM.  
Authors : Zarate-Romero, A.; Rudino-Pinera, E.  
Deposited on : 2013-01-17  
Resolution : 1.95 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](mailto: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>.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	2.36.2
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36.2

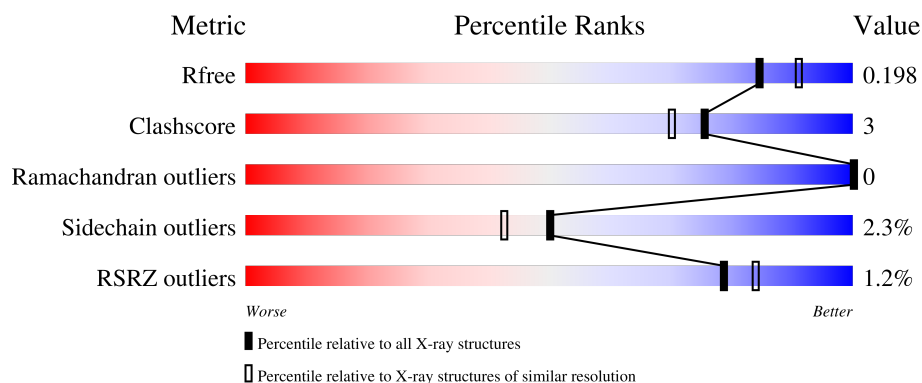
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.95 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))
$R_{free}$	130704	2580 (1.96-1.96)
Clashscore	141614	2705 (1.96-1.96)
Ramachandran outliers	138981	2678 (1.96-1.96)
Sidechain outliers	138945	2678 (1.96-1.96)
RSRZ outliers	127900	2539 (1.96-1.96)

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  $\geq 3$ , 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	746	<div> <div>84%</div> <div>7%</div> <div>9%</div> </div>
1	B	746	<div> <div>83%</div> <div>8%</div> <div>9%</div> </div>
1	C	746	<div> <div>84%</div> <div>7%</div> <div>9%</div> </div>
1	D	746	<div> <div>84%</div> <div>6%</div> <div>9%</div> </div>

## 2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 24049 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	680	Total	C	N	O	S	0	6	0
			5378	3402	952	1018	6			
1	B	677	Total	C	N	O	S	0	9	0
			5379	3403	955	1015	6			
1	C	679	Total	C	N	O	S	0	10	0
			5403	3416	958	1023	6			
1	D	678	Total	C	N	O	S	0	8	0
			5374	3403	950	1015	6			

There are 108 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-26	MET	-	expression tag	UNP Q9C169
A	-25	ASN	-	expression tag	UNP Q9C169
A	-24	HIS	-	expression tag	UNP Q9C169
A	-23	LYS	-	expression tag	UNP Q9C169
A	-22	VAL	-	expression tag	UNP Q9C169
A	-21	HIS	-	expression tag	UNP Q9C169
A	-20	HIS	-	expression tag	UNP Q9C169
A	-19	HIS	-	expression tag	UNP Q9C169
A	-18	HIS	-	expression tag	UNP Q9C169
A	-17	HIS	-	expression tag	UNP Q9C169
A	-16	HIS	-	expression tag	UNP Q9C169
A	-15	ILE	-	expression tag	UNP Q9C169
A	-14	GLU	-	expression tag	UNP Q9C169
A	-13	GLY	-	expression tag	UNP Q9C169
A	-12	ARG	-	expression tag	UNP Q9C169
A	-11	HIS	-	expression tag	UNP Q9C169
A	-10	MET	-	expression tag	UNP Q9C169
A	-9	GLU	-	expression tag	UNP Q9C169
A	-8	LEU	-	expression tag	UNP Q9C169
A	-7	GLY	-	expression tag	UNP Q9C169
A	-6	THR	-	expression tag	UNP Q9C169

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-5	LEU	-	expression tag	UNP Q9C169
A	-4	GLU	-	expression tag	UNP Q9C169
A	-3	GLY	-	expression tag	UNP Q9C169
A	-2	SER	-	expression tag	UNP Q9C169
A	-1	GLU	-	expression tag	UNP Q9C169
A	0	PHE	-	expression tag	UNP Q9C169
B	-26	MET	-	expression tag	UNP Q9C169
B	-25	ASN	-	expression tag	UNP Q9C169
B	-24	HIS	-	expression tag	UNP Q9C169
B	-23	LYS	-	expression tag	UNP Q9C169
B	-22	VAL	-	expression tag	UNP Q9C169
B	-21	HIS	-	expression tag	UNP Q9C169
B	-20	HIS	-	expression tag	UNP Q9C169
B	-19	HIS	-	expression tag	UNP Q9C169
B	-18	HIS	-	expression tag	UNP Q9C169
B	-17	HIS	-	expression tag	UNP Q9C169
B	-16	HIS	-	expression tag	UNP Q9C169
B	-15	ILE	-	expression tag	UNP Q9C169
B	-14	GLU	-	expression tag	UNP Q9C169
B	-13	GLY	-	expression tag	UNP Q9C169
B	-12	ARG	-	expression tag	UNP Q9C169
B	-11	HIS	-	expression tag	UNP Q9C169
B	-10	MET	-	expression tag	UNP Q9C169
B	-9	GLU	-	expression tag	UNP Q9C169
B	-8	LEU	-	expression tag	UNP Q9C169
B	-7	GLY	-	expression tag	UNP Q9C169
B	-6	THR	-	expression tag	UNP Q9C169
B	-5	LEU	-	expression tag	UNP Q9C169
B	-4	GLU	-	expression tag	UNP Q9C169
B	-3	GLY	-	expression tag	UNP Q9C169
B	-2	SER	-	expression tag	UNP Q9C169
B	-1	GLU	-	expression tag	UNP Q9C169
B	0	PHE	-	expression tag	UNP Q9C169
C	-26	MET	-	expression tag	UNP Q9C169
C	-25	ASN	-	expression tag	UNP Q9C169
C	-24	HIS	-	expression tag	UNP Q9C169
C	-23	LYS	-	expression tag	UNP Q9C169
C	-22	VAL	-	expression tag	UNP Q9C169
C	-21	HIS	-	expression tag	UNP Q9C169
C	-20	HIS	-	expression tag	UNP Q9C169
C	-19	HIS	-	expression tag	UNP Q9C169
C	-18	HIS	-	expression tag	UNP Q9C169

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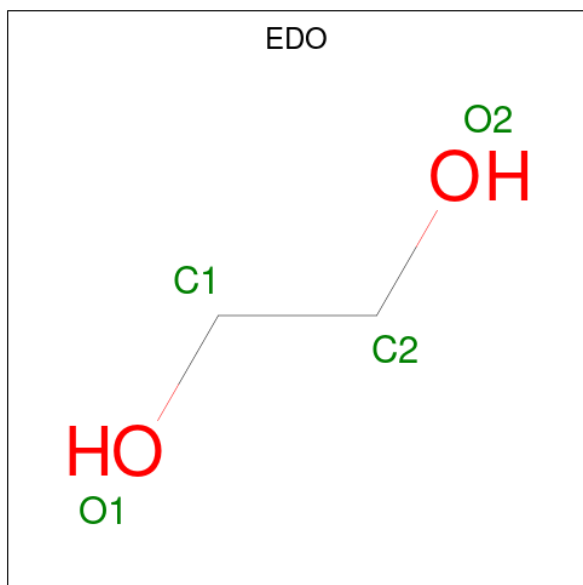
Chain	Residue	Modelled	Actual	Comment	Reference
C	-17	HIS	-	expression tag	UNP Q9C169
C	-16	HIS	-	expression tag	UNP Q9C169
C	-15	ILE	-	expression tag	UNP Q9C169
C	-14	GLU	-	expression tag	UNP Q9C169
C	-13	GLY	-	expression tag	UNP Q9C169
C	-12	ARG	-	expression tag	UNP Q9C169
C	-11	HIS	-	expression tag	UNP Q9C169
C	-10	MET	-	expression tag	UNP Q9C169
C	-9	GLU	-	expression tag	UNP Q9C169
C	-8	LEU	-	expression tag	UNP Q9C169
C	-7	GLY	-	expression tag	UNP Q9C169
C	-6	THR	-	expression tag	UNP Q9C169
C	-5	LEU	-	expression tag	UNP Q9C169
C	-4	GLU	-	expression tag	UNP Q9C169
C	-3	GLY	-	expression tag	UNP Q9C169
C	-2	SER	-	expression tag	UNP Q9C169
C	-1	GLU	-	expression tag	UNP Q9C169
C	0	PHE	-	expression tag	UNP Q9C169
D	-26	MET	-	expression tag	UNP Q9C169
D	-25	ASN	-	expression tag	UNP Q9C169
D	-24	HIS	-	expression tag	UNP Q9C169
D	-23	LYS	-	expression tag	UNP Q9C169
D	-22	VAL	-	expression tag	UNP Q9C169
D	-21	HIS	-	expression tag	UNP Q9C169
D	-20	HIS	-	expression tag	UNP Q9C169
D	-19	HIS	-	expression tag	UNP Q9C169
D	-18	HIS	-	expression tag	UNP Q9C169
D	-17	HIS	-	expression tag	UNP Q9C169
D	-16	HIS	-	expression tag	UNP Q9C169
D	-15	ILE	-	expression tag	UNP Q9C169
D	-14	GLU	-	expression tag	UNP Q9C169
D	-13	GLY	-	expression tag	UNP Q9C169
D	-12	ARG	-	expression tag	UNP Q9C169
D	-11	HIS	-	expression tag	UNP Q9C169
D	-10	MET	-	expression tag	UNP Q9C169
D	-9	GLU	-	expression tag	UNP Q9C169
D	-8	LEU	-	expression tag	UNP Q9C169
D	-7	GLY	-	expression tag	UNP Q9C169
D	-6	THR	-	expression tag	UNP Q9C169
D	-5	LEU	-	expression tag	UNP Q9C169
D	-4	GLU	-	expression tag	UNP Q9C169
D	-3	GLY	-	expression tag	UNP Q9C169

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-2	SER	-	expression tag	UNP Q9C169
D	-1	GLU	-	expression tag	UNP Q9C169
D	0	PHE	-	expression tag	UNP Q9C169

- Molecule 2 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C<sub>2</sub>H<sub>6</sub>O<sub>2</sub>).



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

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

- Molecule 3 is TETRAETHYLENE GLYCOL (three-letter code: PG4) (formula:  $C_8H_{18}O_5$ ).



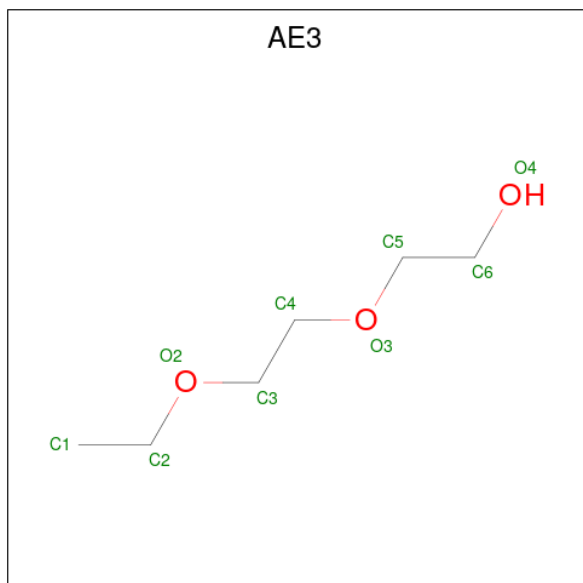
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			6	4	2		
3	A	1	Total	C	O	0	0
			6	4	2		
3	B	1	Total	C	O	0	0
			6	4	2		
3	B	1	Total	C	O	0	0
			6	4	2		
3	B	1	Total	C	O	0	0
			6	4	2		
3	C	1	Total	C	O	0	0
			6	4	2		
3	C	1	Total	C	O	0	0
			6	4	2		
3	C	1	Total	C	O	0	0
			6	4	2		
3	C	1	Total	C	O	0	0
			12	8	4		
3	C	1	Total	C	O	0	0
			12	8	4		

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

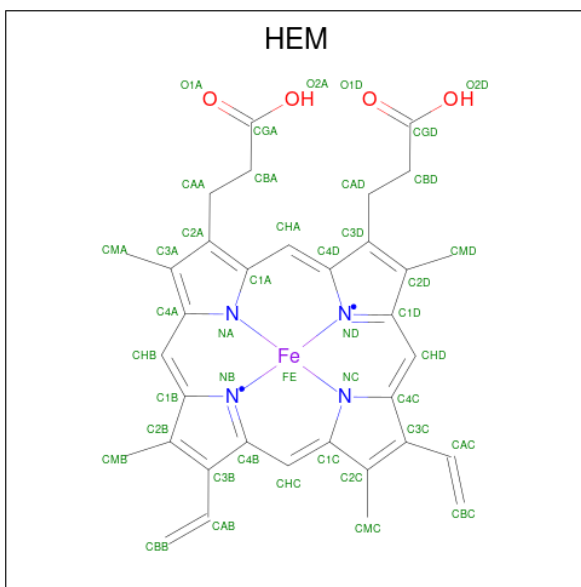
- Molecule 4 is 2-(2-ETHOXYETHOXY)ETHANOL (three-letter code: AE3) (formula:  $C_6H_{14}O_3$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			9	6	3		
4	B	1	Total	C	O	0	0
			9	6	3		
4	D	1	Total	C	O	0	0
			9	6	3		

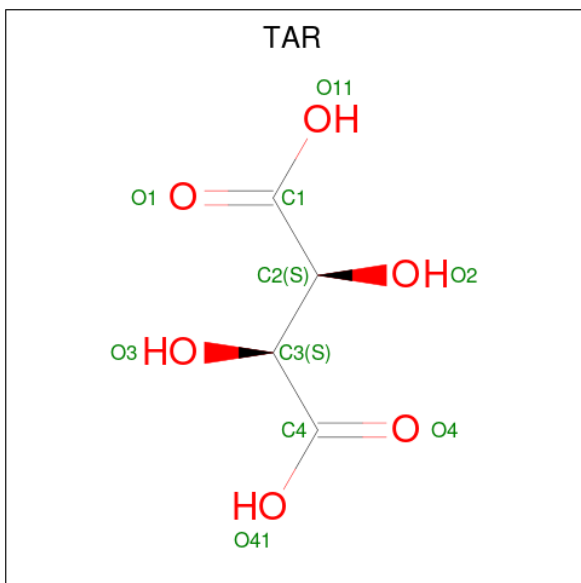
- Molecule 5 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (formula:  $C_{34}H_{32}FeN_4O_4$ ).





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

- Molecule 6 is D(-)-TARTARIC ACID (three-letter code: TAR) (formula:  $C_4H_6O_6$ ).



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

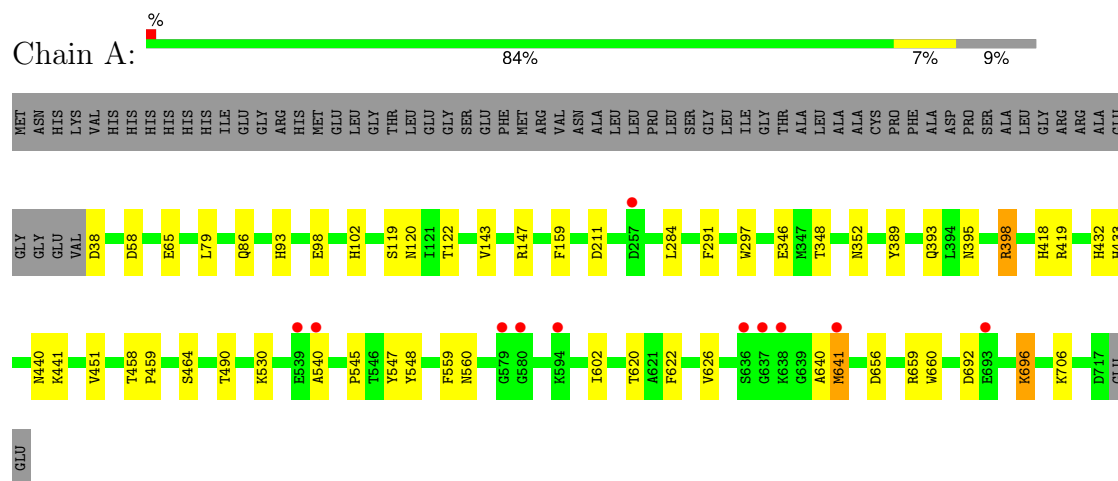
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	633	Total	O	0	0
			633	633		
7	B	611	Total	O	0	0
			611	611		
7	C	501	Total	O	0	0
			501	501		
7	D	444	Total	O	0	0
			444	444		

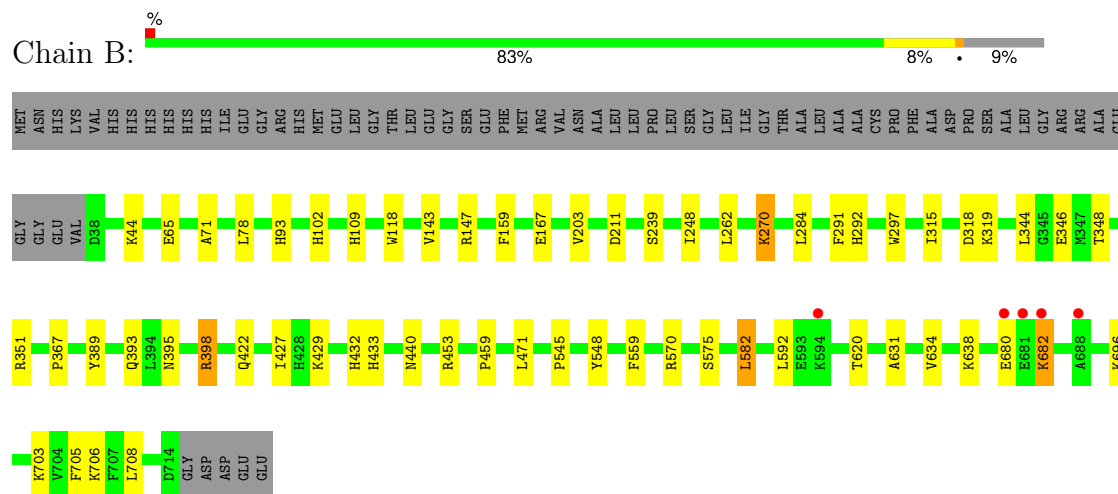
### 3 Residue-property plots

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.

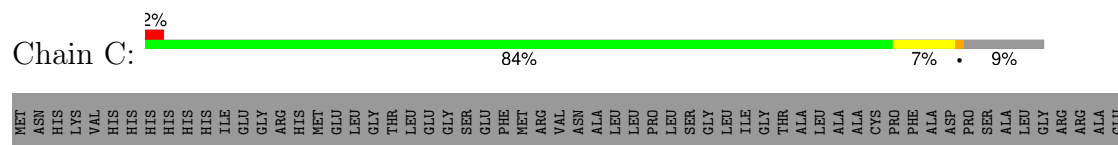
#### • Molecule 1: CATALASE-3

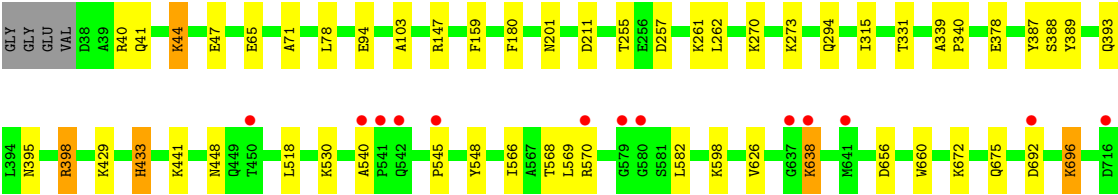


#### • Molecule 1: CATALASE-3



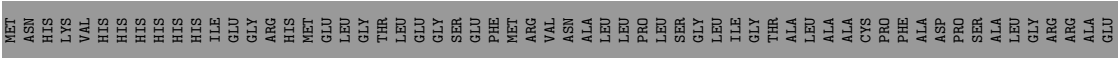
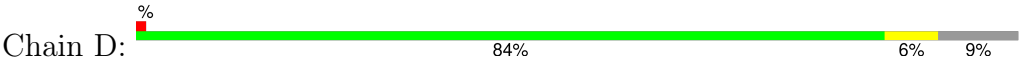
#### • Molecule 1: CATALASE-3





ASP  
GLU  
GLU

• Molecule 1: CATALASE-3



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	131.52Å 154.80Å 160.74Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	42.52 – 1.95 42.52 – 1.95	Depositor EDS
% Data completeness (in resolution range)	94.7 (42.52-1.95) 94.7 (42.52-1.95)	Depositor EDS
$R_{merge}$	0.14	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.33 (at 1.95Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, $R_{free}$	0.157 , 0.197 0.158 , 0.198	Depositor DCC
$R_{free}$ test set	11250 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	12.1	Xtriage
Anisotropy	0.092	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 50.4	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.015 for -h,l,k	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	24049	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	11.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.57% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: AE3, PG4, HEM, TAR, EDO

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	0.38	0/5511	0.53	0/7475
1	B	0.42	0/5513	0.55	0/7476
1	C	0.38	0/5536	0.55	0/7508
1	D	0.38	0/5508	0.53	0/7469
All	All	0.39	0/22068	0.54	0/29928

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	C	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	568	THR	Mainchain

### 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	5378	0	5193	40	0
1	B	5379	0	5199	42	0
1	C	5403	0	5218	36	0
1	D	5374	0	5196	32	0
2	A	6	0	6	0	0
2	B	15	0	15	0	0
2	C	3	0	3	0	0
2	D	9	0	9	1	0
3	A	12	0	14	3	0
3	B	18	0	21	0	0
3	C	42	0	51	0	0
3	D	12	0	14	1	0
4	A	9	0	14	0	0
4	B	9	0	14	0	0
4	D	9	0	14	0	0
5	A	43	0	30	0	0
5	B	43	0	30	1	0
5	C	43	0	30	0	0
5	D	43	0	30	1	0
6	C	10	0	4	2	0
7	A	633	0	0	9	0
7	B	611	0	0	6	0
7	C	501	0	0	5	0
7	D	444	0	0	5	0
All	All	24049	0	21105	135	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 (135) 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)
7:A:2230:HOH:O	1:D:147[B]:ARG:O	1.82	0.95
1:B:147[A]:ARG:O	7:B:2113:HOH:O	1.85	0.94
7:B:2223:HOH:O	1:C:147[A]:ARG:O	1.90	0.90
1:C:40:ARG:NH2	1:C:47:GLU:OE2	2.10	0.84
1:A:147[B]:ARG:O	7:A:2114:HOH:O	2.05	0.74
1:A:441:LYS:HG2	3:A:1721:PG4:H21	1.74	0.70
1:A:640:ALA:O	7:A:2599:HOH:O	2.10	0.68
1:B:570[B]:ARG:HG2	1:B:570[B]:ARG:HH11	1.58	0.68
1:B:398:ARG:HD3	1:D:422:GLN:NE2	2.09	0.67
1:C:273:LYS:HZ3	6:C:1717:TAR:H3	1.61	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:38:ASP:N	7:A:2001:HOH:O	2.29	0.64
1:C:638:LYS:HE2	1:C:638:LYS:H	1.62	0.64
1:B:422:GLN:NE2	1:D:398:ARG:HG2	2.14	0.63
5:D:4000:HEM:HBB2	5:D:4000:HEM:HMB2	1.81	0.62
1:C:273:LYS:NZ	6:C:1717:TAR:H3	2.15	0.61
1:C:675:GLN:OE1	7:C:2467:HOH:O	2.17	0.60
1:B:248:ILE:HD12	5:B:4000:HEM:HMB1	1.84	0.59
1:D:346:GLU:OE2	1:D:348:THR:OG1	2.14	0.59
1:D:395:ASN:OD1	1:D:398:ARG:NH1	2.35	0.58
1:B:398:ARG:HG2	1:C:398:ARG:NH1	2.18	0.58
7:C:2294:HOH:O	1:D:398:ARG:NH1	2.35	0.58
1:A:398:ARG:NH2	1:B:422:GLN:HE21	2.02	0.58
1:A:398:ARG:NE	1:D:398:ARG:HE	2.03	0.55
1:B:459[A]:PRO:O	7:B:2268:HOH:O	2.18	0.54
1:A:419:ARG:O	1:C:94:GLU:HG2	2.08	0.54
1:A:398:ARG:HE	1:D:398:ARG:HE	1.55	0.53
1:B:346:GLU:OE2	1:B:348:THR:OG1	2.15	0.53
1:B:570[B]:ARG:HH11	1:B:570[B]:ARG:CG	2.22	0.52
1:D:578:LYS:HG3	1:D:633:ARG:CZ	2.40	0.52
1:A:441:LYS:HZ2	3:A:1721:PG4:H12	1.76	0.51
1:C:71:ALA:HB2	1:C:78:LEU:HD21	1.92	0.51
1:A:147[B]:ARG:HD2	7:A:2155:HOH:O	2.10	0.50
1:C:570[B]:ARG:HA	1:C:598:LYS:O	2.12	0.50
1:D:631:ALA:O	1:D:634:VAL:HG22	2.11	0.50
1:A:395:ASN:O	1:A:398:ARG:HD2	2.12	0.49
1:C:545:PRO:HA	1:C:548:TYR:CD2	2.48	0.49
1:D:559:PHE:HE1	1:D:706:LYS:HG3	1.77	0.49
1:B:389:TYR:O	1:B:393:GLN:HG2	2.12	0.49
1:C:530:LYS:HG2	1:C:540:ALA:HB1	1.93	0.49
1:B:147[B]:ARG:HD2	7:B:2157:HOH:O	2.12	0.49
1:C:103:ALA:O	1:C:147[B]:ARG:HG2	2.14	0.48
1:A:284:LEU:HD13	1:A:620:THR:HB	1.95	0.48
1:B:147[A]:ARG:HD2	7:B:2157:HOH:O	2.12	0.48
1:A:58:ASP:HB2	7:A:2028:HOH:O	2.12	0.48
1:A:346:GLU:OE2	1:A:348:THR:OG1	2.21	0.48
1:C:638:LYS:H	1:C:638:LYS:CE	2.27	0.47
1:A:102:HIS:CE1	1:A:143:VAL:HG22	2.48	0.47
1:C:378:GLU:N	1:C:378:GLU:OE1	2.48	0.47
1:A:432:HIS:HB2	1:A:440:ASN:HB3	1.96	0.47
1:C:692:ASP:O	1:C:696:LYS:HD3	2.14	0.47
1:C:429:LYS:HB2	1:C:429:LYS:HE3	1.71	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:395:ASN:O	1:B:398:ARG:HD2	2.14	0.46
1:C:570[A]:ARG:HA	1:C:598:LYS:O	2.15	0.46
1:B:318:ASP:OD1	1:B:319:LYS:HG3	2.15	0.46
1:D:656:ASP:HB3	1:D:660:TRP:CZ3	2.50	0.46
1:B:203:VAL:HG11	1:C:294:GLN:NE2	2.30	0.46
1:B:682:LYS:H	1:B:682:LYS:NZ	2.13	0.46
1:A:119:SER:HA	1:A:122:THR:O	2.16	0.46
1:C:262:LEU:HB2	1:C:315:ILE:HB	1.97	0.46
1:C:270:LYS:HA	1:C:270:LYS:HE2	1.95	0.46
1:A:93:HIS:ND1	1:D:395:ASN:HB2	2.30	0.46
1:B:71:ALA:HB2	1:B:78:LEU:HD21	1.98	0.46
1:B:102:HIS:CE1	1:B:143:VAL:HG22	2.50	0.46
1:B:429:LYS:HE3	1:B:429:LYS:HB2	1.55	0.45
1:A:696:LYS:HB3	1:A:696:LYS:HE3	1.46	0.45
1:B:270:LYS:NZ	7:B:2281:HOH:O	2.48	0.45
1:A:389:TYR:O	1:A:393:GLN:HG2	2.17	0.45
1:B:559:PHE:HB3	1:B:708:LEU:HD11	1.99	0.45
1:D:429:LYS:HE3	1:D:429:LYS:HB2	1.76	0.45
1:A:490:THR:HG21	1:A:560:ASN:HB3	1.98	0.45
1:A:545:PRO:HA	1:A:548:TYR:CD2	2.52	0.45
1:D:297:TRP:CZ3	1:D:352:ASN:HB3	2.51	0.45
1:B:545:PRO:HA	1:B:548:TYR:CD2	2.53	0.44
1:C:433:HIS:ND1	1:C:433:HIS:N	2.64	0.44
1:A:86:GLN:HB2	7:A:2028:HOH:O	2.17	0.44
1:B:284:LEU:HD21	1:B:291:PHE:CD2	2.53	0.44
1:D:474:THR:HA	3:D:1719:PG4:H21	2.00	0.44
1:D:593:GLU:OE2	7:D:2391:HOH:O	2.21	0.44
1:A:297:TRP:CZ3	1:A:352:ASN:HB3	2.53	0.44
1:A:297:TRP:CE3	1:D:201:ASN:HB3	2.53	0.44
1:B:118:TRP:CZ2	1:B:344:LEU:HA	2.53	0.44
1:C:389:TYR:O	1:C:393:GLN:HG2	2.17	0.44
1:A:98:GLU:HB3	1:A:147[B]:ARG:HG3	2.00	0.44
1:C:41:GLN:HA	1:C:44:LYS:HE3	2.00	0.44
1:A:490:THR:CG2	1:A:560:ASN:HB3	2.48	0.43
1:B:575:SER:HA	1:B:631:ALA:HB2	2.00	0.43
1:A:418:HIS:C	1:A:419:ARG:HG2	2.39	0.43
1:B:631:ALA:O	1:B:634:VAL:HG22	2.18	0.43
1:C:180:PHE:O	1:C:331:THR:HG22	2.18	0.43
1:C:255:THR:OG1	1:C:257:ASP:OD1	2.24	0.43
1:B:427:ILE:O	1:D:47:GLU:HA	2.18	0.43
1:D:432:HIS:HB2	1:D:440:ASN:HB3	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:389:TYR:O	1:D:393:GLN:HG2	2.18	0.43
1:C:570[A]:ARG:HB3	7:C:2428:HOH:O	2.17	0.43
1:B:432:HIS:HB2	1:B:440:ASN:HB3	2.00	0.43
1:B:682:LYS:HE3	1:B:682:LYS:HB3	1.60	0.43
1:C:656:ASP:HB3	1:C:660:TRP:CZ3	2.54	0.43
1:A:398:ARG:HH21	1:B:422:GLN:HE21	1.63	0.42
1:A:602:ILE:HD11	1:A:622:PHE:CZ	2.54	0.42
1:B:284:LEU:HD13	1:B:620:THR:HB	1.99	0.42
1:B:705:PHE:CD2	1:B:706:LYS:HG2	2.53	0.42
1:B:93:HIS:ND1	1:C:395:ASN:HB2	2.33	0.42
1:C:261:LYS:HE2	7:C:2246:HOH:O	2.19	0.42
1:D:476:ASP:O	7:D:2314:HOH:O	2.21	0.42
1:B:297:TRP:CE3	1:C:201:ASN:HB3	2.54	0.42
1:D:591:GLN:O	1:D:594:LYS:HG2	2.19	0.42
1:A:559:PHE:HE1	1:A:706:LYS:HG3	1.84	0.42
1:A:547:TYR:CE1	3:A:1720:PG4:H21	2.55	0.42
1:D:343[A]:VAL:HG22	7:D:2247:HOH:O	2.19	0.42
1:A:284:LEU:HD21	1:A:291:PHE:CD2	2.55	0.42
7:A:2046:HOH:O	1:C:448:ASN:HB2	2.19	0.42
1:D:570:ARG:NE	7:D:2151:HOH:O	2.41	0.42
1:D:409:ARG:HH12	2:D:1716:EDO:H21	1.85	0.42
1:C:566:ILE:HB	1:C:569[B]:LEU:HD12	2.00	0.41
1:B:395:ASN:O	1:B:398:ARG:NH1	2.53	0.41
1:C:339:ALA:HA	1:C:340:PRO:HD2	1.88	0.41
1:D:545:PRO:HA	1:D:548:TYR:CD2	2.55	0.41
1:D:554:ARG:NH1	7:D:2242:HOH:O	2.33	0.41
1:B:582:LEU:HD12	1:B:582:LEU:HA	1.89	0.41
1:D:494:GLN:O	1:D:498:ILE:HG12	2.21	0.41
1:A:530:LYS:HG2	1:A:540:ALA:HB1	2.03	0.41
1:B:109:HIS:HB2	1:B:351:ARG:HB2	2.03	0.41
1:A:79:LEU:HD11	1:C:387:TYR:HB2	2.02	0.41
1:B:167:GLU:OE2	1:B:453:ARG:NE	2.46	0.41
1:B:262:LEU:HB2	1:B:315:ILE:HB	2.03	0.41
1:B:367:PRO:HD3	1:B:389:TYR:CD2	2.56	0.41
1:A:458:THR:HA	1:A:459:PRO:HD3	1.95	0.41
1:D:262:LEU:HB2	1:D:315:ILE:HG12	2.03	0.41
1:A:641:MET:HE2	1:A:641:MET:HB3	1.88	0.40
1:A:656:ASP:HB3	1:A:660:TRP:CZ3	2.56	0.40
1:A:395:ASN:HB2	1:D:93:HIS:ND1	2.36	0.40
1:A:659:ARG:HD2	7:A:2612:HOH:O	2.21	0.40
1:C:441:LYS:NZ	7:C:2334:HOH:O	2.35	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:239:SER:HA	1:B:292:HIS:CD2	2.56	0.40
1:D:261:LYS:HB3	1:D:261:LYS:HE3	1.88	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	684/746 (92%)	664 (97%)	20 (3%)	0	100	100
1	B	684/746 (92%)	659 (96%)	25 (4%)	0	100	100
1	C	687/746 (92%)	668 (97%)	19 (3%)	0	100	100
1	D	684/746 (92%)	662 (97%)	22 (3%)	0	100	100
All	All	2739/2984 (92%)	2653 (97%)	86 (3%)	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	563/609 (92%)	549 (98%)	14 (2%)	47	38
1	B	563/609 (92%)	548 (97%)	15 (3%)	44	34
1	C	566/609 (93%)	551 (97%)	15 (3%)	44	34

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	D	563/609 (92%)	551 (98%)	12 (2%)	53	46
All	All	2255/2436 (93%)	2199 (98%)	56 (2%)	50	38

All (56) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	65	GLU
1	A	120	ASN
1	A	159	PHE
1	A	211	ASP
1	A	398	ARG
1	A	433	HIS
1	A	451[A]	VAL
1	A	451[B]	VAL
1	A	464	SER
1	A	626[A]	VAL
1	A	626[B]	VAL
1	A	641	MET
1	A	692	ASP
1	A	696	LYS
1	B	44	LYS
1	B	65	GLU
1	B	159	PHE
1	B	211	ASP
1	B	270	LYS
1	B	398	ARG
1	B	433	HIS
1	B	471	LEU
1	B	582	LEU
1	B	592	LEU
1	B	638	LYS
1	B	680	GLU
1	B	682	LYS
1	B	696	LYS
1	B	703	LYS
1	C	44	LYS
1	C	65	GLU
1	C	159	PHE
1	C	211	ASP
1	C	388[A]	SER
1	C	388[B]	SER
1	C	398	ARG

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Mol	Chain	Res	Type
1	C	433	HIS
1	C	518[A]	LEU
1	C	518[B]	LEU
1	C	582	LEU
1	C	626	VAL
1	C	638	LYS
1	C	672	LYS
1	C	696	LYS
1	D	65	GLU
1	D	159	PHE
1	D	211	ASP
1	D	294	GLN
1	D	315	ILE
1	D	429	LYS
1	D	433	HIS
1	D	560	ASN
1	D	582	LEU
1	D	608	SER
1	D	659	ARG
1	D	692	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	422	GLN

### 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 monosaccharides in this entry.

## 5.6 Ligand geometry

31 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$
3	PG4	A	1720	-	5,5,12	0.52	0	4,4,11	1.66	1 (25%)
2	EDO	B	1715	-	2,2,3	0.36	0	1,1,2	0.26	0
2	EDO	C	1718	-	2,2,3	0.37	0	1,1,2	0.28	0
3	PG4	B	1721	-	5,5,12	0.52	0	4,4,11	1.24	1 (25%)
5	HEM	A	4000	1	42,50,50	1.90	6 (14%)	46,82,82	1.41	5 (10%)
2	EDO	B	1717	-	2,2,3	0.34	0	1,1,2	0.47	0
3	PG4	C	1720	-	5,5,12	0.49	0	4,4,11	1.52	0
3	PG4	C	1721	-	5,5,12	0.54	0	4,4,11	1.41	0
2	EDO	D	1717	-	2,2,3	0.34	0	1,1,2	0.25	0
4	AE3	D	1721	-	8,8,8	0.57	0	7,7,7	1.76	2 (28%)
2	EDO	B	1719	-	2,2,3	0.33	0	1,1,2	0.23	0
3	PG4	C	1719	-	5,5,12	0.54	0	4,4,11	1.45	1 (25%)
5	HEM	B	4000	1	42,50,50	1.90	8 (19%)	46,82,82	1.47	4 (8%)
2	EDO	B	1716	-	2,2,3	0.34	0	1,1,2	0.37	0
4	AE3	B	1723	-	8,8,8	0.62	0	7,7,7	1.44	0
5	HEM	D	4000	1	42,50,50	1.91	8 (19%)	46,82,82	1.52	4 (8%)
4	AE3	A	1722	-	8,8,8	0.55	0	7,7,7	1.47	0
3	PG4	B	1720	-	5,5,12	0.55	0	4,4,11	1.43	0
2	EDO	D	1718	-	2,2,3	0.34	0	1,1,2	0.32	0
5	HEM	C	4000	1	42,50,50	1.88	7 (16%)	46,82,82	1.47	7 (15%)
2	EDO	A	1719	-	2,2,3	0.35	0	1,1,2	0.31	0
3	PG4	C	1723	-	11,11,12	0.89	0	10,10,11	1.56	1 (10%)
3	PG4	B	1722	-	5,5,12	0.51	0	4,4,11	1.55	0
2	EDO	D	1716	-	2,2,3	0.36	0	1,1,2	0.35	0
2	EDO	A	1718	-	2,2,3	0.36	0	1,1,2	0.44	0
3	PG4	A	1721	-	5,5,12	0.51	0	4,4,11	1.55	1 (25%)
6	TAR	C	1717	-	9,9,9	1.60	2 (22%)	12,12,12	1.44	2 (16%)
3	PG4	D	1720	-	5,5,12	0.50	0	4,4,11	1.46	1 (25%)
3	PG4	D	1719	-	5,5,12	0.54	0	4,4,11	1.46	1 (25%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
3	PG4	C	1722	-	11,11,12	0.91	0	10,10,11	1.44	0
2	EDO	B	1718	-	2,2,3	0.36	0	1,1,2	0.17	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
3	PG4	A	1720	-	-	2/3/3/10	-
3	PG4	B	1721	-	-	2/3/3/10	-
5	HEM	A	4000	1	-	2/12/54/54	-
3	PG4	C	1720	-	-	2/3/3/10	-
3	PG4	C	1721	-	-	1/3/3/10	-
4	AE3	D	1721	-	-	4/6/6/6	-
3	PG4	C	1719	-	-	0/3/3/10	-
5	HEM	D	4000	1	-	4/12/54/54	-
4	AE3	B	1723	-	-	2/6/6/6	-
5	HEM	B	4000	1	-	2/12/54/54	-
4	AE3	A	1722	-	-	3/6/6/6	-
3	PG4	B	1720	-	-	3/3/3/10	-
5	HEM	C	4000	1	-	2/12/54/54	-
3	PG4	C	1723	-	-	2/9/9/10	-
3	PG4	B	1722	-	-	0/3/3/10	-
3	PG4	A	1721	-	-	1/3/3/10	-
6	TAR	C	1717	-	-	8/12/12/12	-
3	PG4	D	1720	-	-	2/3/3/10	-
3	PG4	D	1719	-	-	1/3/3/10	-
3	PG4	C	1722	-	-	5/9/9/10	-

All (31) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	4000	HEM	C3D-C2D	7.66	1.53	1.36
5	D	4000	HEM	C3D-C2D	7.59	1.53	1.36
5	C	4000	HEM	C3D-C2D	7.37	1.52	1.36
5	B	4000	HEM	C3D-C2D	7.30	1.52	1.36
5	A	4000	HEM	C3C-C2C	-4.28	1.34	1.40
5	C	4000	HEM	C3C-C2C	-4.18	1.34	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	4000	HEM	C3C-C2C	-4.07	1.34	1.40
5	D	4000	HEM	C3C-C2C	-3.64	1.35	1.40
5	C	4000	HEM	C3C-CAC	3.42	1.55	1.47
5	D	4000	HEM	C3C-CAC	3.39	1.55	1.47
5	B	4000	HEM	C3C-CAC	3.37	1.55	1.47
5	B	4000	HEM	CAB-C3B	3.22	1.56	1.47
5	A	4000	HEM	CAB-C3B	3.17	1.55	1.47
5	D	4000	HEM	CAB-C3B	3.15	1.55	1.47
5	A	4000	HEM	C3C-CAC	3.09	1.54	1.47
5	C	4000	HEM	CAB-C3B	2.82	1.54	1.47
5	D	4000	HEM	FE-ND	2.73	2.13	1.98
5	D	4000	HEM	CMB-C2B	2.56	1.56	1.50
5	B	4000	HEM	FE-ND	2.48	2.11	1.98
5	B	4000	HEM	C3C-C4C	2.47	1.45	1.41
5	C	4000	HEM	C3C-C4C	2.43	1.44	1.41
5	A	4000	HEM	C3C-C4C	2.37	1.44	1.41
5	C	4000	HEM	FE-ND	2.36	2.11	1.98
5	A	4000	HEM	CMB-C2B	2.33	1.55	1.50
5	D	4000	HEM	C3C-C4C	2.30	1.44	1.41
5	B	4000	HEM	CMB-C2B	2.27	1.55	1.50
5	C	4000	HEM	CMB-C2B	2.25	1.55	1.50
5	D	4000	HEM	CMD-C2D	2.23	1.55	1.50
6	C	1717	TAR	O2-C2	-2.21	1.37	1.42
6	C	1717	TAR	O3-C3	-2.08	1.38	1.42
5	B	4000	HEM	CMD-C2D	2.01	1.54	1.50

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	D	4000	HEM	C4D-ND-C1D	5.16	111.32	105.21
5	C	4000	HEM	C4D-ND-C1D	5.02	111.15	105.21
5	B	4000	HEM	C4D-ND-C1D	4.97	111.10	105.21
5	A	4000	HEM	C4D-ND-C1D	4.73	110.80	105.21
5	D	4000	HEM	CAA-CBA-CGA	-3.43	104.58	113.83
5	A	4000	HEM	CAA-CBA-CGA	-3.35	104.80	113.83
5	B	4000	HEM	CAA-CBA-CGA	-3.22	105.17	113.83
5	C	4000	HEM	C4C-CHD-C1D	3.17	126.75	122.56
6	C	1717	TAR	O11-C1-C2	3.02	121.71	113.31
5	C	4000	HEM	CAA-CBA-CGA	-2.85	106.16	113.83
5	B	4000	HEM	C4C-CHD-C1D	2.63	126.03	122.56
5	C	4000	HEM	CBD-CAD-C3D	-2.56	105.45	112.53
5	D	4000	HEM	C4C-CHD-C1D	2.56	125.94	122.56

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	4000	HEM	CBD-CAD-C3D	-2.40	105.90	112.53
4	D	1721	AE3	O2-C3-C4	2.39	121.23	110.35
3	A	1720	PG4	C3-O2-C2	2.35	121.31	113.06
5	A	4000	HEM	CAD-C3D-C4D	2.27	128.65	124.70
5	D	4000	HEM	CBD-CAD-C3D	-2.24	106.34	112.53
5	A	4000	HEM	C4C-CHD-C1D	2.21	125.48	122.56
3	C	1719	PG4	C3-O2-C2	2.17	120.69	113.06
5	A	4000	HEM	CMA-C3A-C4A	-2.11	125.36	128.46
5	C	4000	HEM	C1B-NB-C4B	2.11	107.70	105.21
3	B	1721	PG4	C3-O2-C2	2.08	120.37	113.06
3	A	1721	PG4	C3-O2-C2	2.05	120.25	113.06
3	D	1719	PG4	O2-C2-C1	2.05	119.13	110.11
6	C	1717	TAR	O41-C4-O4	-2.05	119.44	124.08
4	D	1721	AE3	C5-O3-C4	2.04	122.17	113.26
3	C	1723	PG4	C7-O4-C6	2.04	120.21	113.06
5	C	4000	HEM	CMD-C2D-C1D	2.03	128.21	125.03
3	D	1720	PG4	C3-O2-C2	2.02	120.13	113.06
5	C	4000	HEM	C3B-C4B-NB	-2.01	108.03	109.47

There are no chirality outliers.

All (48) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	C	1717	TAR	O3-C3-C4-O4
6	C	1717	TAR	O2-C2-C3-O3
3	C	1722	PG4	C5-C6-O4-C7
6	C	1717	TAR	C1-C2-C3-C4
6	C	1717	TAR	O3-C3-C4-O41
4	A	1722	AE3	O2-C3-C4-O3
6	C	1717	TAR	C1-C2-C3-O3
6	C	1717	TAR	O2-C2-C3-C4
4	D	1721	AE3	O2-C3-C4-O3
6	C	1717	TAR	O1-C1-C2-O2
6	C	1717	TAR	O11-C1-C2-O2
4	D	1721	AE3	C6-C5-O3-C4
3	B	1720	PG4	C1-C2-O2-C3
3	B	1721	PG4	O1-C1-C2-O2
3	C	1722	PG4	O3-C5-C6-O4
3	B	1720	PG4	O1-C1-C2-O2
3	D	1720	PG4	C4-C3-O2-C2
4	A	1722	AE3	C6-C5-O3-C4
3	C	1720	PG4	O1-C1-C2-O2

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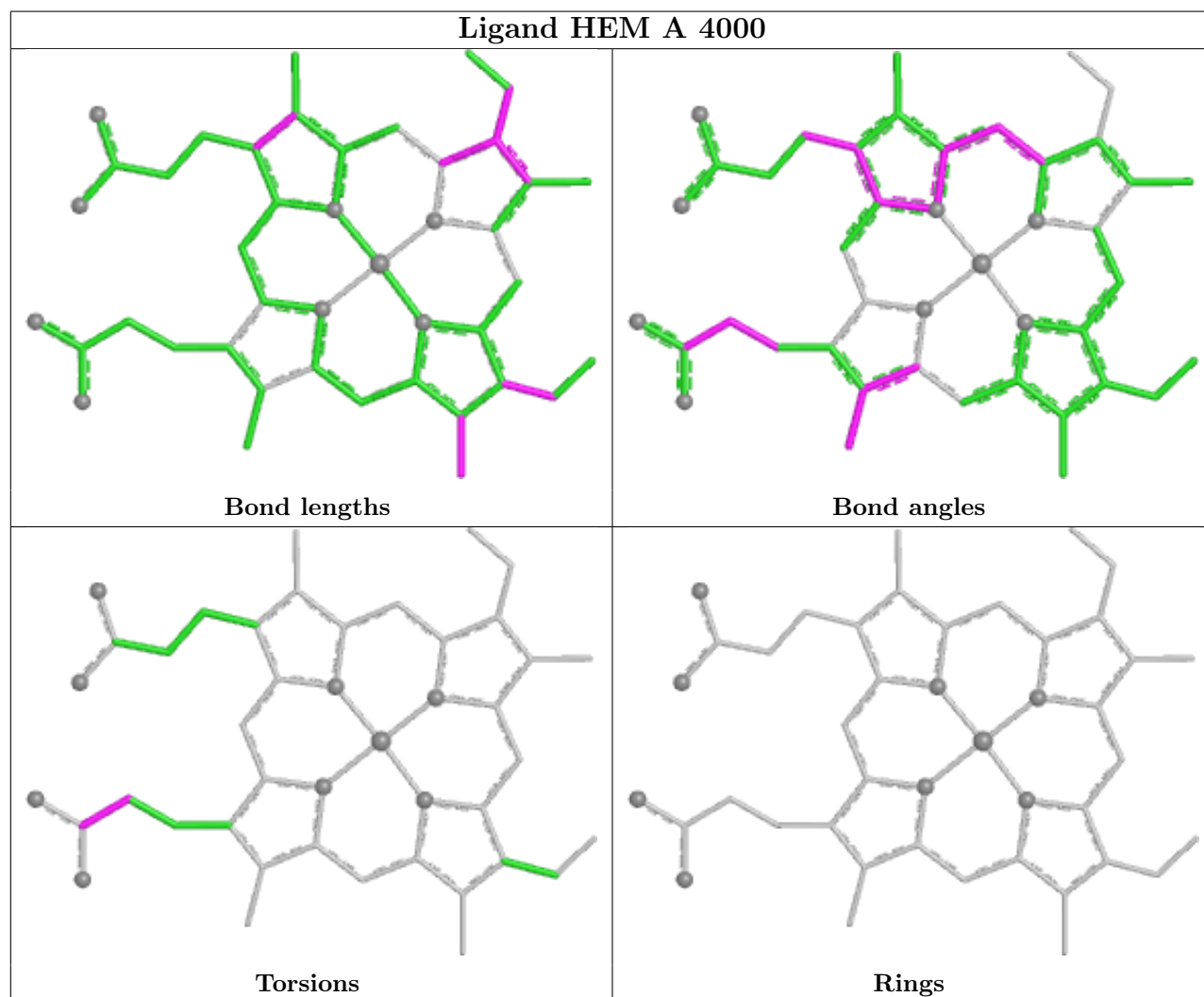
Mol	Chain	Res	Type	Atoms
4	A	1722	AE3	C1-C2-O2-C3
3	B	1720	PG4	C4-C3-O2-C2
4	B	1723	AE3	C1-C2-O2-C3
3	C	1720	PG4	C4-C3-O2-C2
4	B	1723	AE3	C3-C4-O3-C5
4	D	1721	AE3	O3-C5-C6-O4
3	D	1720	PG4	C1-C2-O2-C3
3	C	1722	PG4	C3-C4-O3-C5
3	C	1722	PG4	C4-C3-O2-C2
3	D	1719	PG4	C4-C3-O2-C2
3	C	1723	PG4	C5-C6-O4-C7
5	D	4000	HEM	CAA-CBA-CGA-O2A
5	A	4000	HEM	CAA-CBA-CGA-O2A
5	C	4000	HEM	CAA-CBA-CGA-O1A
4	D	1721	AE3	C4-C3-O2-C2
5	B	4000	HEM	CAA-CBA-CGA-O1A
5	C	4000	HEM	CAA-CBA-CGA-O2A
5	B	4000	HEM	CAA-CBA-CGA-O2A
5	D	4000	HEM	CAA-CBA-CGA-O1A
5	A	4000	HEM	CAA-CBA-CGA-O1A
3	A	1720	PG4	C1-C2-O2-C3
3	C	1721	PG4	C4-C3-O2-C2
3	B	1721	PG4	C4-C3-O2-C2
3	C	1722	PG4	C8-C7-O4-C6
3	A	1720	PG4	O1-C1-C2-O2
3	C	1723	PG4	O3-C5-C6-O4
3	A	1721	PG4	C1-C2-O2-C3
5	D	4000	HEM	CAD-CBD-CGD-O1D
5	D	4000	HEM	CAD-CBD-CGD-O2D

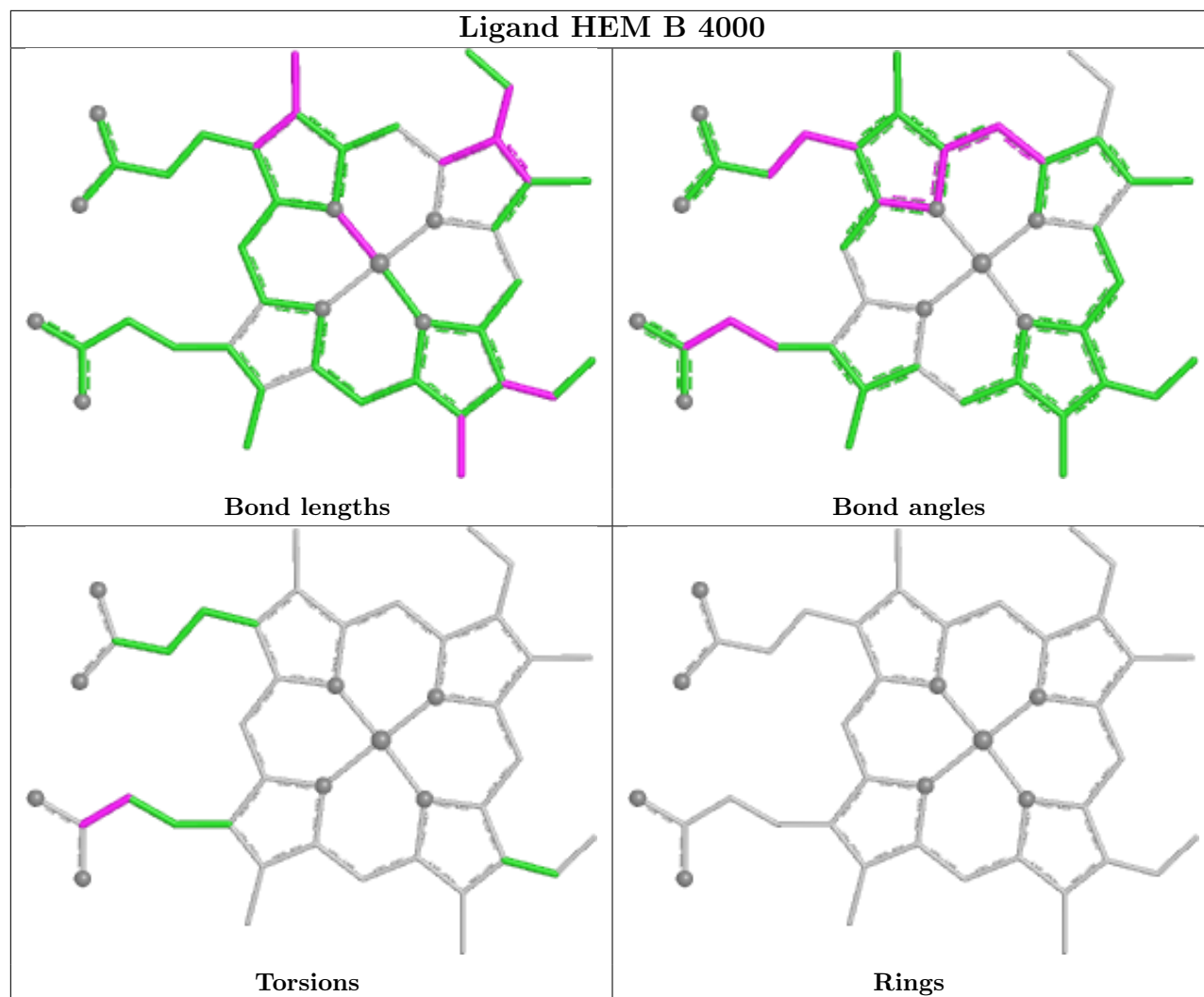
There are no ring outliers.

7 monomers are involved in 9 short contacts:

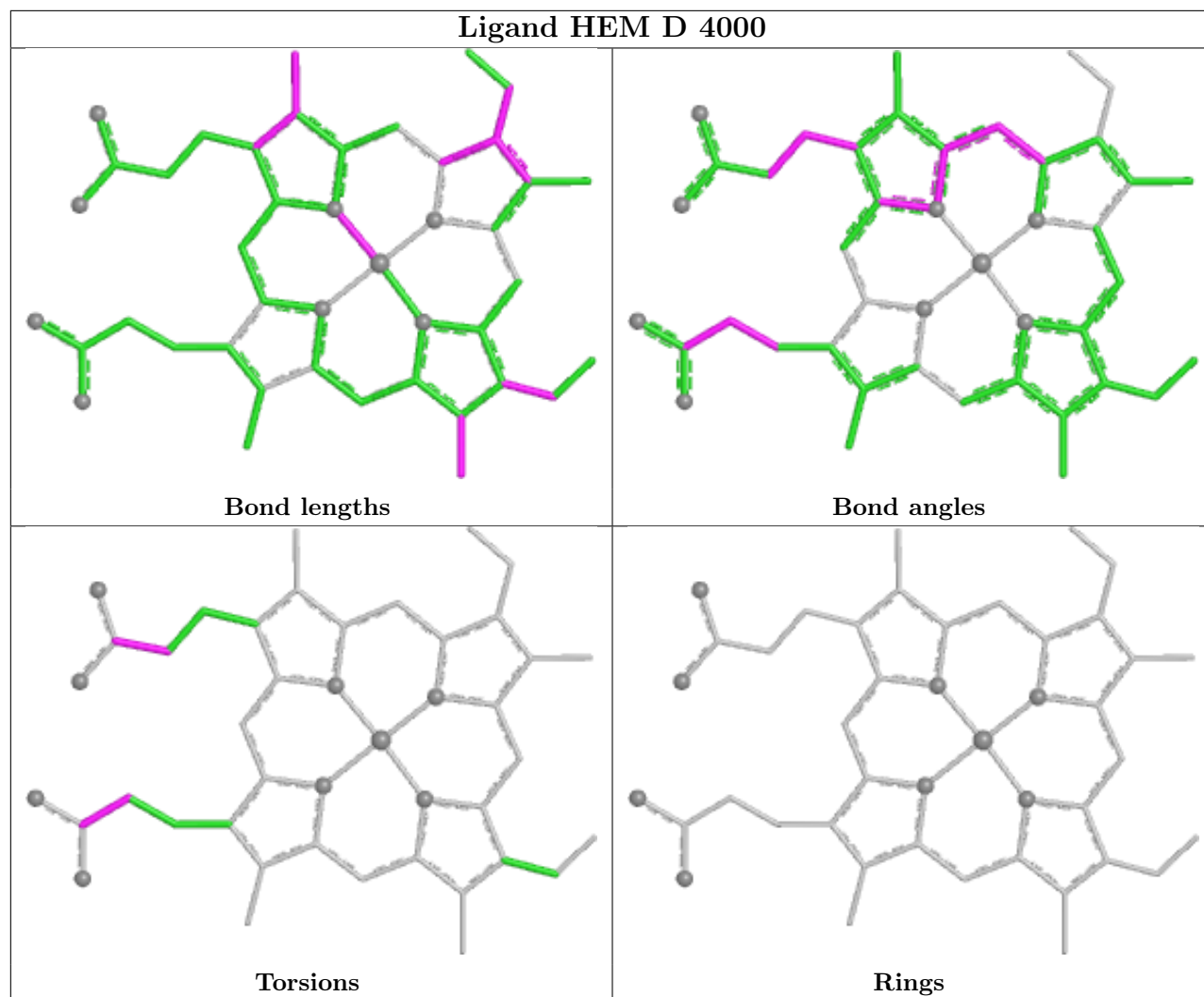
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1720	PG4	1	0
5	B	4000	HEM	1	0
5	D	4000	HEM	1	0
2	D	1716	EDO	1	0
3	A	1721	PG4	2	0
6	C	1717	TAR	2	0
3	D	1719	PG4	1	0

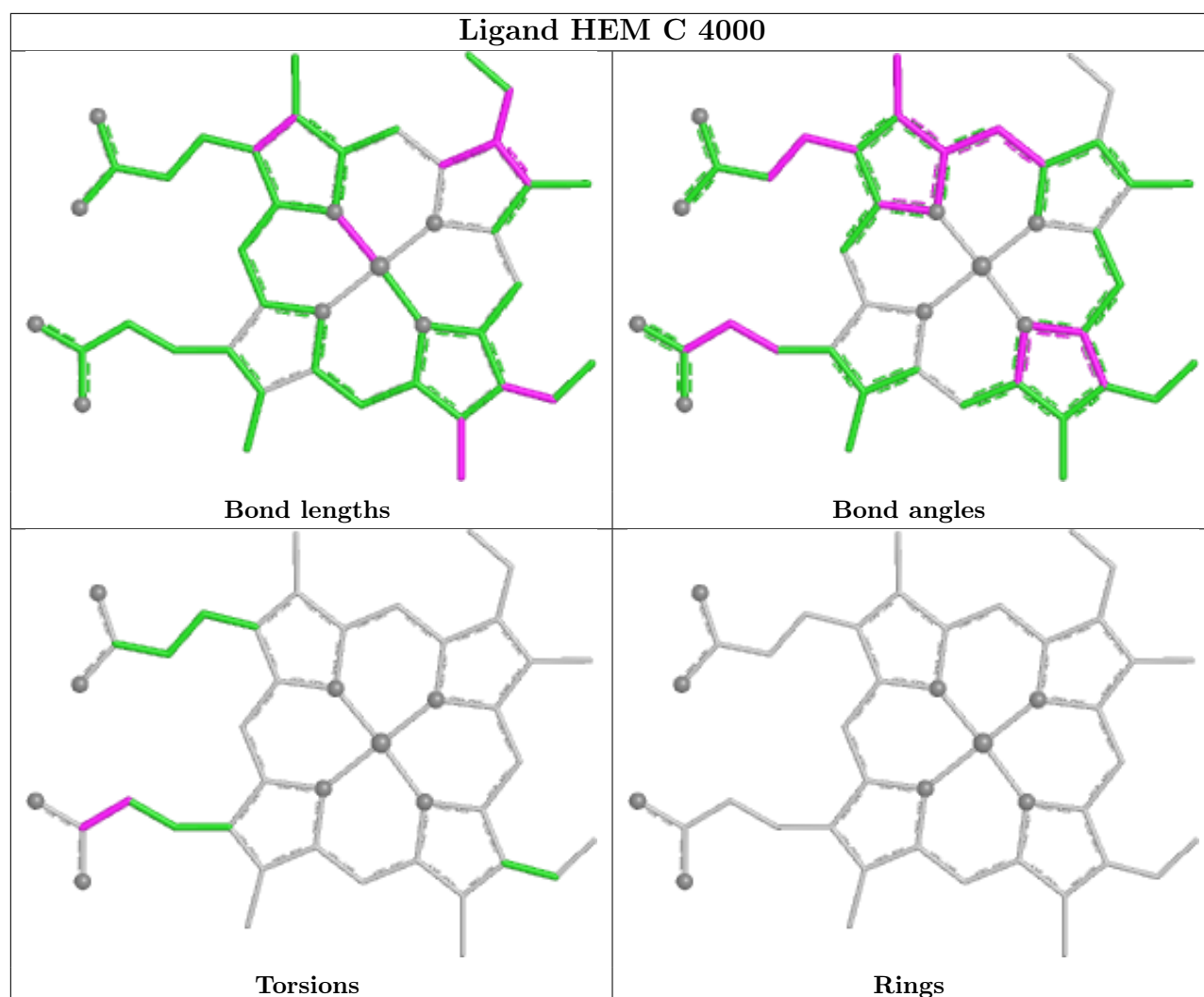
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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.





## Ligand HEM D 4000





## 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 ⓘ

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	680/746 (91%)	-0.15	11 (1%) 72 79	1, 10, 28, 61	0
1	B	677/746 (90%)	-0.21	5 (0%) 87 92	1, 7, 26, 50	0
1	C	679/746 (91%)	-0.20	13 (1%) 66 74	1, 8, 26, 64	0
1	D	678/746 (90%)	-0.19	4 (0%) 89 93	1, 8, 26, 54	0
All	All	2714/2984 (90%)	-0.19	33 (1%) 79 84	1, 8, 26, 64	0

All (33) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	540	ALA	4.3
1	C	540	ALA	4.2
1	A	641	MET	4.2
1	C	641	MET	3.5
1	A	638	LYS	3.5
1	C	638	LYS	3.0
1	A	637	GLY	3.0
1	C	637	GLY	2.9
1	B	682	LYS	2.9
1	A	636	SER	2.8
1	C	570[A]	ARG	2.6
1	C	450	THR	2.5
1	B	594	LYS	2.4
1	A	579	GLY	2.4
1	D	594	LYS	2.4
1	D	682	LYS	2.4
1	C	580	GLY	2.4
1	C	541	PRO	2.4
1	A	539	GLU	2.3
1	B	680	GLU	2.3
1	C	579	GLY	2.3

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Mol	Chain	Res	Type	RSRZ
1	C	545	PRO	2.3
1	C	716	ASP	2.3
1	C	542	GLN	2.2
1	D	638	LYS	2.2
1	D	692	ASP	2.2
1	B	681	GLU	2.1
1	C	692	ASP	2.1
1	A	580	GLY	2.1
1	A	693	GLU	2.0
1	A	594	LYS	2.0
1	B	688	ALA	2.0
1	A	257	ASP	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 monosaccharides in this entry.

## 6.4 Ligands [i](#)

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, 95<sup>th</sup> 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( $\text{\AA}^2$ )	Q<0.9
2	EDO	B	1718	3/4	0.70	0.20	21,21,38,43	0
3	PG4	B	1721	6/13	0.70	0.24	36,43,46,55	0
4	AE3	D	1721	9/9	0.70	0.26	35,43,66,66	0
4	AE3	A	1722	9/9	0.74	0.24	32,39,52,65	0
2	EDO	A	1719	3/4	0.79	0.15	24,24,30,36	0
3	PG4	C	1723	12/13	0.80	0.22	16,21,28,31	0
2	EDO	D	1717	3/4	0.81	0.16	27,27,28,37	0
2	EDO	C	1718	3/4	0.82	0.15	29,29,38,43	0
2	EDO	B	1715	3/4	0.82	0.13	25,25,33,35	0
3	PG4	B	1722	6/13	0.84	0.21	24,25,37,42	0
4	AE3	B	1723	9/9	0.84	0.19	26,36,50,54	0

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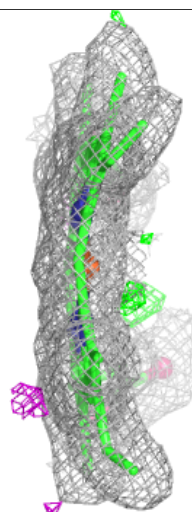
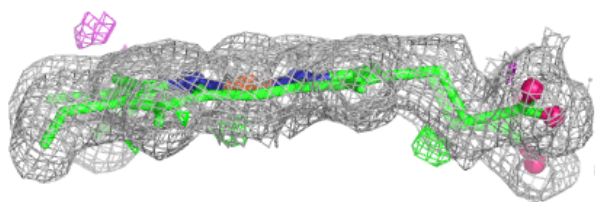
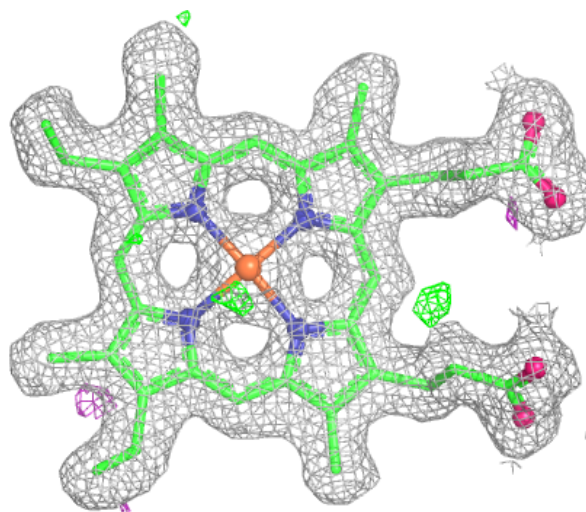
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	EDO	B	1717	3/4	0.84	0.27	26,26,40,45	0
2	EDO	D	1716	3/4	0.85	0.14	23,23,27,36	0
3	PG4	A	1721	6/13	0.86	0.14	18,29,32,37	0
3	PG4	D	1719	6/13	0.87	0.14	22,29,37,37	0
3	PG4	B	1720	6/13	0.88	0.16	25,33,40,40	0
2	EDO	B	1716	3/4	0.88	0.12	24,24,24,36	0
3	PG4	C	1722	12/13	0.89	0.12	20,28,37,50	0
2	EDO	A	1718	3/4	0.89	0.12	28,28,31,37	0
3	PG4	C	1720	6/13	0.89	0.12	28,31,37,39	0
6	TAR	C	1717	10/10	0.89	0.17	10,26,35,39	0
2	EDO	B	1719	3/4	0.90	0.14	19,19,24,30	0
3	PG4	C	1719	6/13	0.91	0.18	11,31,36,38	0
3	PG4	D	1720	6/13	0.91	0.20	25,31,37,51	0
3	PG4	C	1721	6/13	0.91	0.13	22,26,37,39	0
2	EDO	D	1718	3/4	0.92	0.08	19,19,21,30	0
3	PG4	A	1720	6/13	0.93	0.10	13,22,23,32	0
5	HEM	B	4000	43/43	0.98	0.09	0,2,6,8	0
5	HEM	C	4000	43/43	0.98	0.10	0,2,4,8	0
5	HEM	A	4000	43/43	0.98	0.10	0,4,7,7	0
5	HEM	D	4000	43/43	0.99	0.10	0,2,5,8	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.

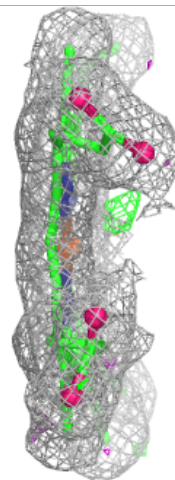
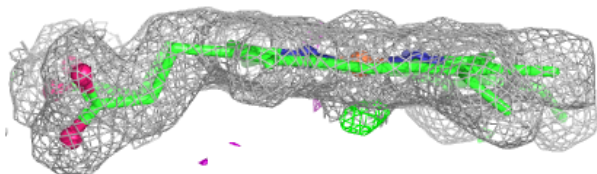
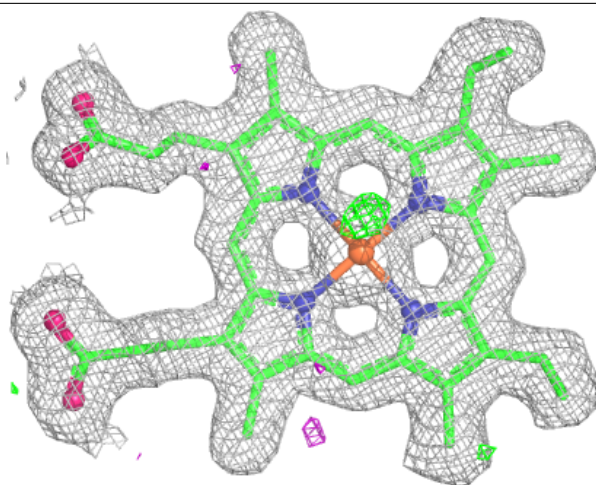
**Electron density around HEM B 4000:**

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



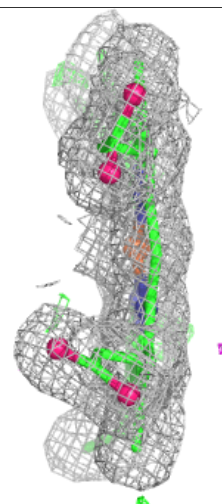
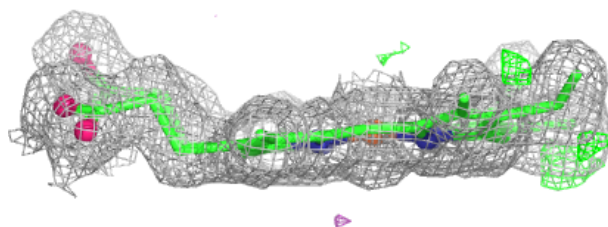
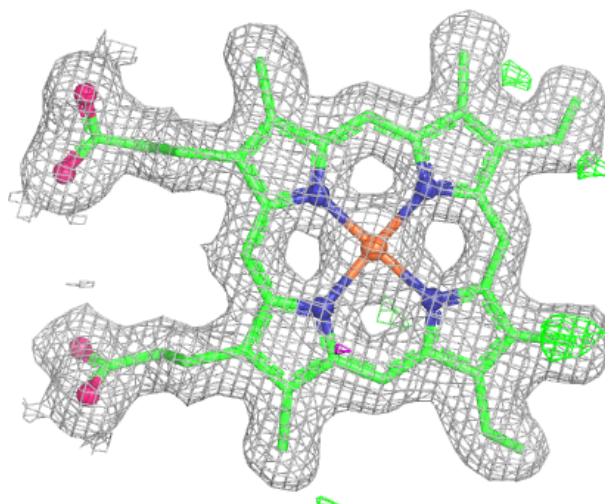
**Electron density around HEM C 4000:**

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



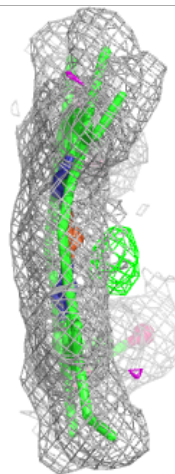
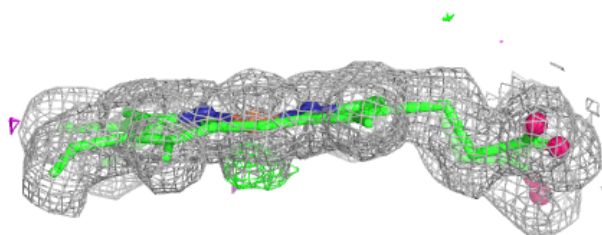
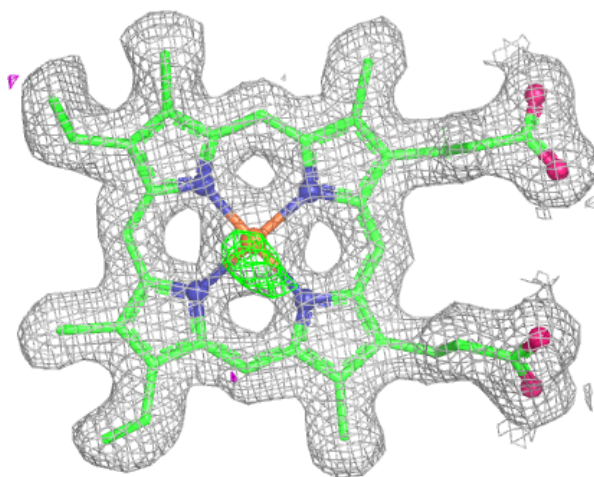
**Electron density around HEM A 4000:**

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



**Electron density around HEM D 4000:**

$2mF_o-DF_c$  (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.