



# Full wwPDB X-ray Structure Validation Report ⓘ

Apr 20, 2021 – 09:01 AM JST

PDB ID : 6LFK  
Title : Crystal structure of KatE from atypical E. coli  
Authors : Cho, H.S.; Yoo, Y.; Park, J.B.  
Deposited on : 2019-12-03  
Resolution : 2.10 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.

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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 : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.18  
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) : 2.18

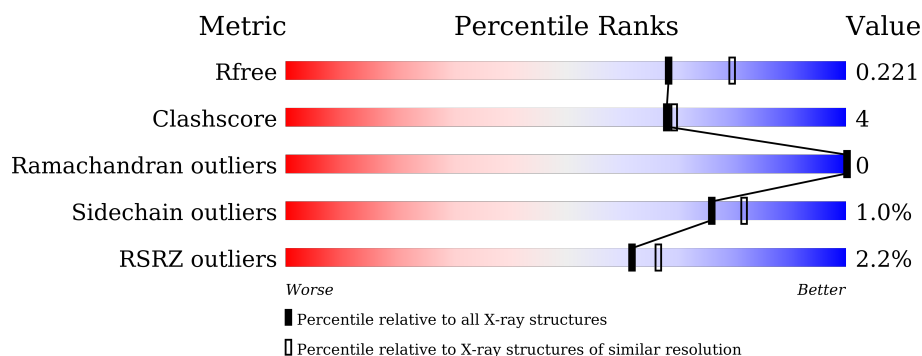
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.10 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	5197 (2.10-2.10)
Clashscore	141614	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (2.10-2.10)
RSRZ outliers	127900	5083 (2.10-2.10)

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	750	<div> <div></div> <div>85%11%.</div> </div>
1	B	750	<div> <div></div> <div>85%12%.</div> </div>
1	C	750	<div> <div>2%</div> <div>87%9%.</div> </div>
1	D	750	<div> <div>3%</div> <div>86%10%.</div> </div>

## 2 Entry composition [i](#)

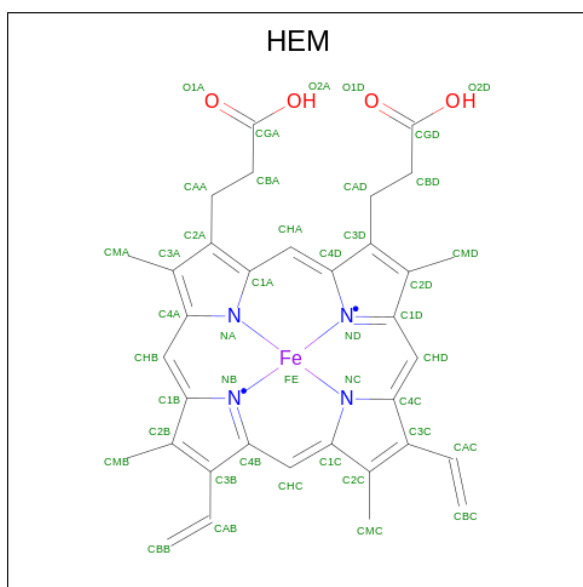
There are 3 unique types of molecules in this entry. The entry contains 23791 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 eKatE catalase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	719	Total	C	N	O	S	0	0	0
			5725	3633	1002	1069	21			
1	B	721	Total	C	N	O	S	0	0	0
			5737	3641	1004	1071	21			
1	C	723	Total	C	N	O	S	0	0	0
			5755	3652	1007	1075	21			
1	D	720	Total	C	N	O	S	0	0	0
			5730	3636	1003	1070	21			

- Molecule 2 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: 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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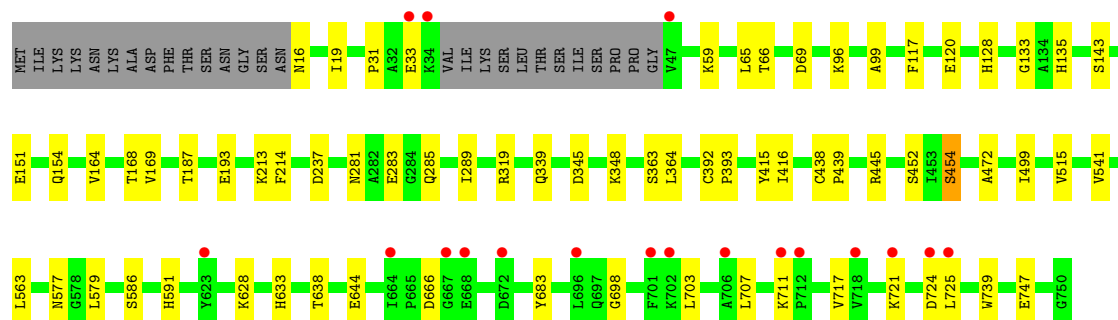
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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	C	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
2	D	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		

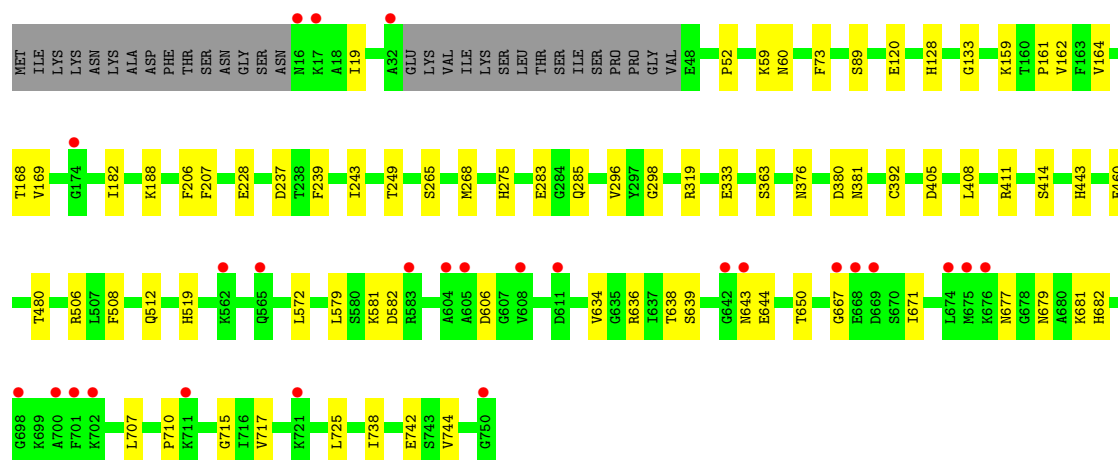
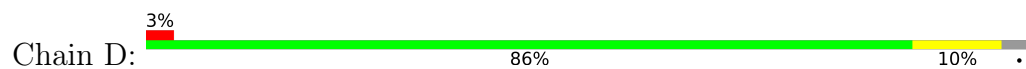
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	149	Total	O	0	0
			149	149		
3	B	196	Total	O	0	0
			196	196		
3	C	166	Total	O	0	0
			166	166		
3	D	161	Total	O	0	0
			161	161		





• Molecule 1: eKatE catalase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	76.58Å 178.94Å 107.28Å 90.00° 106.70° 90.00°	Depositor
Resolution (Å)	47.49 – 2.10 47.49 – 2.10	Depositor EDS
% Data completeness (in resolution range)	97.1 (47.49-2.10) 97.1 (47.49-2.10)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.80 (at 2.10Å)	Xtriage
Refinement program	PHENIX 1.14_3260	Depositor
R, $R_{free}$	0.188 , 0.223 0.186 , 0.221	Depositor DCC
$R_{free}$ test set	1262 reflections (0.81%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	31.8	Xtriage
Anisotropy	0.382	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 42.1	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	23791	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.82% 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: 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	0.47	0/5877	0.60	0/7970
1	B	0.49	0/5890	0.61	0/7989
1	C	0.47	0/5908	0.59	0/8012
1	D	0.48	0/5883	0.59	0/7979
All	All	0.48	0/23558	0.60	0/31950

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	5725	0	5565	52	0
1	B	5737	0	5577	53	0
1	C	5755	0	5596	45	0
1	D	5730	0	5568	45	0
2	A	43	0	30	1	0
2	B	43	0	30	3	0
2	C	43	0	30	1	0
2	D	43	0	30	2	0
3	A	149	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	196	0	0	0	0
3	C	166	0	0	1	0
3	D	161	0	0	0	0
All	All	23791	0	22426	174	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 4.

All (174) 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:C:392:CYS:SG	1:C:415:TYR:HB2	1.13	1.65
1:C:392:CYS:SG	1:C:415:TYR:CB	2.05	1.44
1:A:452:SER:HB3	1:B:73:PHE:HE2	1.32	0.94
1:C:666:ASP:OD2	1:C:698:GLY:HA3	1.78	0.84
1:D:710:PRO:HG3	1:D:744:VAL:HG21	1.76	0.67
1:A:647:VAL:HG12	1:A:649:GLY:H	1.62	0.65
1:D:363:SER:HG	1:D:582:ASP:CG	2.00	0.64
1:C:392:CYS:SG	1:C:415:TYR:HB3	2.31	0.63
1:B:474:PRO:HG3	1:B:480:THR:HG21	1.80	0.63
1:A:128:HIS:CE1	1:A:169:VAL:HG22	2.35	0.62
1:B:214:PHE:CD2	2:B:801:HEM:HAC	2.34	0.62
1:A:452:SER:HB3	1:B:73:PHE:CE2	2.24	0.61
1:A:685:ILE:HG13	1:A:707:LEU:HD13	1.82	0.61
1:D:606:ASP:OD1	1:D:638:THR:N	2.33	0.58
1:B:142:GLU:HA	1:B:156:PRO:HG3	1.85	0.57
1:C:717:VAL:HG12	1:C:725:LEU:HD12	1.87	0.56
1:B:229:ILE:HD11	1:D:319:ARG:HG2	1.88	0.56
1:A:363:SER:HB2	1:A:579:LEU:HD11	1.88	0.55
1:C:33:GLU:O	1:C:33:GLU:HG2	2.07	0.55
1:C:452:SER:HB2	1:D:73:PHE:CE2	2.42	0.54
1:A:454:SER:O	1:B:381:ASN:HB2	2.08	0.54
1:B:178:THR:HG21	1:B:310:LEU:HD23	1.89	0.54
1:D:443:HIS:CE1	1:D:460:GLU:HB2	2.42	0.54
1:B:443:HIS:CE1	1:B:460:GLU:HB2	2.42	0.54
1:A:64:LYS:HE2	1:A:190:TYR:CE1	2.42	0.53
1:C:638:THR:HG22	1:C:644:GLU:HG2	1.90	0.53
1:D:636:ARG:HD3	1:D:644:GLU:HG2	1.90	0.53
1:D:363:SER:OG	1:D:582:ASP:OD2	2.09	0.52
1:A:535:TRP:CD2	1:C:31:PRO:HG3	2.44	0.52
1:D:275:HIS:CG	1:D:408:LEU:HB2	2.44	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:579:LEU:HD22	1:A:581:LYS:O	2.10	0.52
1:A:567:GLN:HA	1:A:570:HIS:CE1	2.45	0.51
1:D:639:SER:OG	1:D:643:ASN:N	2.36	0.51
1:B:362:GLU:HG2	1:B:367:VAL:HG23	1.92	0.51
1:D:681:LYS:HB3	1:D:707:LEU:HD21	1.92	0.51
1:A:614:ASP:O	1:A:618:LYS:HG2	2.11	0.51
1:B:664:ILE:HD12	1:B:696:LEU:HD23	1.92	0.50
1:B:689:LYS:NZ	1:B:745:ALA:O	2.44	0.50
1:B:499:ILE:HG21	1:D:19:ILE:HG12	1.93	0.50
1:D:634:VAL:HG13	1:D:650:THR:HG22	1.94	0.50
1:B:267:ARG:HG2	1:B:332:PRO:HG3	1.92	0.50
1:C:281:ASN:ND2	1:C:285:GLN:HB2	2.27	0.50
1:D:363:SER:HB2	1:D:579:LEU:HD21	1.94	0.50
1:B:531:VAL:O	1:B:537:ARG:NH1	2.43	0.50
1:A:415:TYR:O	1:A:419:GLN:HG2	2.12	0.49
1:C:193:GLU:OE2	1:C:193:GLU:HA	2.12	0.49
1:D:677:ASN:OD1	1:D:679:ASN:HB2	2.11	0.49
1:A:712:PRO:HB3	1:A:716:ILE:HB	1.93	0.49
1:B:128:HIS:HA	1:B:168:THR:O	2.13	0.49
1:C:289:ILE:HA	1:C:339:GLN:O	2.12	0.49
1:B:639:SER:OG	1:B:641:GLN:O	2.22	0.48
1:B:420:ILE:HD12	1:B:426:VAL:HG12	1.96	0.48
1:C:454:SER:O	1:D:381:ASN:HB2	2.14	0.48
1:B:473:PRO:HD2	1:B:476:GLU:HG3	1.96	0.47
1:D:405:ASP:HB3	1:D:408:LEU:HB3	1.97	0.47
1:B:515:VAL:HG21	1:B:739:TRP:HB3	1.95	0.47
1:A:445:ARG:O	1:B:120:GLU:HG2	2.14	0.47
1:A:711:LYS:HE3	1:A:712:PRO:HD2	1.97	0.47
1:A:65:LEU:HD13	1:A:377:ARG:HD3	1.96	0.47
1:B:211:ALA:CB	1:B:410:GLY:HA3	2.45	0.47
1:B:749:ALA:HB2	1:D:681:LYS:HE3	1.95	0.47
1:C:66:THR:HA	1:C:69:ASP:OD2	2.15	0.47
1:C:96:LYS:NZ	1:C:99:ALA:O	2.43	0.47
1:B:474:PRO:HG3	1:B:480:THR:CG2	2.43	0.47
1:B:622:LYS:HE2	1:B:623:TYR:CZ	2.49	0.47
1:C:345:ASP:OD1	1:C:348:LYS:NZ	2.48	0.47
1:D:283:GLU:HB3	1:D:285:GLN:HG3	1.97	0.47
1:A:211:ALA:CB	1:A:410:GLY:HA3	2.45	0.46
1:A:281:ASN:HD21	1:A:285:GLN:HB2	1.80	0.46
1:C:392:CYS:SG	1:C:415:TYR:CG	2.99	0.46
1:C:143:SER:HB2	1:C:154:GLN:O	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:717:VAL:HG12	1:D:725:LEU:HD12	1.96	0.46
1:B:667:GLY:O	1:B:671:ILE:HG12	2.15	0.46
1:B:221:VAL:HG23	1:B:239:PHE:CD1	2.50	0.46
1:B:641:GLN:O	1:B:643:ASN:N	2.45	0.46
1:C:472:ALA:O	1:D:89:SER:HA	2.16	0.46
1:C:515:VAL:HG21	1:C:739:TRP:HB3	1.97	0.46
1:C:133:GLY:HA2	1:C:164:VAL:O	2.16	0.45
1:A:381:ASN:HB2	1:B:454:SER:O	2.15	0.45
1:A:281:ASN:ND2	1:A:285:GLN:HB2	2.31	0.45
1:A:19:ILE:HG12	1:C:499:ILE:HG21	1.98	0.45
1:A:143:SER:HB2	1:A:154:GLN:O	2.16	0.45
1:D:207:PHE:O	1:D:249:THR:HA	2.17	0.45
1:D:667:GLY:O	1:D:671:ILE:HG12	2.16	0.45
1:A:228:GLU:HB3	1:C:319:ARG:NH2	2.31	0.45
1:A:533:ARG:HD3	1:A:535:TRP:CH2	2.52	0.45
1:A:715:GLY:HA3	1:A:738:ILE:HG13	1.98	0.45
1:C:16:ASN:ND2	1:C:19:ILE:HG13	2.32	0.45
1:B:136:GLY:HA3	1:B:374:VAL:O	2.17	0.44
1:D:508:PHE:O	1:D:512:GLN:HG2	2.17	0.44
1:B:713:ASP:OD1	1:B:716:ILE:HG12	2.17	0.44
1:A:128:HIS:HA	1:A:168:THR:O	2.18	0.44
1:A:496:SER:HA	1:A:499:ILE:HD12	1.99	0.44
1:B:133:GLY:HA2	1:B:164:VAL:O	2.16	0.44
1:B:289:ILE:HA	1:B:339:GLN:O	2.17	0.44
1:D:506:ARG:NH1	1:D:572:LEU:HD13	2.33	0.44
1:D:52:PRO:HB3	1:D:380:ASP:HA	2.00	0.44
1:B:647:VAL:HG12	1:B:649:GLY:H	1.82	0.44
1:C:703:LEU:O	1:C:707:LEU:HG	2.18	0.44
1:C:586:SER:O	1:C:591:HIS:HD2	2.01	0.43
1:D:128:HIS:HA	1:D:168:THR:O	2.18	0.43
1:D:715:GLY:HA3	1:D:738:ILE:HG13	1.98	0.43
1:A:689:LYS:HZ1	1:A:750:GLY:H	1.66	0.43
1:A:631:ALA:O	1:A:650:THR:HA	2.18	0.43
1:B:512:GLN:HB2	1:B:517:GLN:HG3	2.00	0.43
1:D:128:HIS:CE1	1:D:169:VAL:HG22	2.53	0.43
1:A:52:PRO:HB3	1:A:380:ASP:HA	1.99	0.43
1:A:229:ILE:HD11	1:C:319:ARG:HG2	1.99	0.43
1:B:333:GLU:HG2	1:B:374:VAL:HG22	2.00	0.43
1:B:528:LEU:HA	1:B:528:LEU:HD23	1.81	0.43
1:A:162:VAL:HA	1:A:188:LYS:O	2.19	0.43
1:A:444:GLN:O	1:A:445:ARG:HD3	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:50:MET:O	1:B:51:MET:HG2	2.18	0.43
1:A:601:ALA:HB2	1:A:659:VAL:HG11	1.99	0.43
1:B:508:PHE:O	1:B:512:GLN:HG2	2.19	0.43
1:A:221:VAL:HG23	1:A:239:PHE:CD1	2.53	0.43
1:A:64:LYS:HE2	1:A:190:TYR:CZ	2.54	0.42
1:A:164:VAL:HG22	1:A:187:THR:HG23	2.01	0.42
1:D:581:LYS:HE3	1:D:581:LYS:HB2	1.84	0.42
1:B:612:ALA:O	1:B:616:ILE:HG13	2.18	0.42
1:C:445:ARG:O	1:D:120:GLU:HG2	2.20	0.42
1:A:124:GLU:HB2	1:A:173:ARG:HG3	2.02	0.42
1:C:363:SER:OG	1:C:579:LEU:HD21	2.19	0.42
1:C:721:LYS:HB3	1:C:721:LYS:HE3	1.81	0.42
1:D:206:PHE:CG	2:D:801:HEM:HAB	2.54	0.42
1:A:558:ASN:O	1:C:633:HIS:HB2	2.20	0.42
1:A:576:ILE:O	1:A:579:LEU:HB2	2.20	0.42
1:C:128:HIS:HA	1:C:168:THR:O	2.20	0.42
1:B:241:ASP:HB2	1:B:536:ILE:HD13	2.00	0.42
1:A:131:GLY:O	1:A:319:ARG:NH2	2.48	0.42
1:B:144:LEU:HD23	1:B:144:LEU:HA	1.82	0.42
1:B:214:PHE:CD2	2:B:801:HEM:CAC	3.02	0.42
1:B:239:PHE:O	1:B:243:ILE:HG12	2.19	0.42
1:C:65:LEU:HD21	1:C:135:HIS:CG	2.55	0.42
1:A:48:GLU:OE2	1:A:56:LYS:NZ	2.53	0.41
1:B:267:ARG:NH2	1:B:321:GLU:OE2	2.35	0.41
1:C:117:PHE:HA	1:C:120:GLU:HG3	2.01	0.41
1:C:164:VAL:HG22	1:C:187:THR:HG23	2.02	0.41
1:A:634:VAL:HG13	1:A:650:THR:HG23	2.02	0.41
1:A:676:LYS:O	1:C:747:GLU:HG2	2.20	0.41
1:A:702:LYS:HE3	1:A:702:LYS:HB2	1.32	0.41
1:B:206:PHE:CG	2:B:801:HEM:HAB	2.55	0.41
1:D:162:VAL:HA	1:D:188:LYS:O	2.20	0.41
1:A:92:GLN:HA	1:C:213:LYS:HD3	2.02	0.41
1:A:603:LEU:O	1:A:665:PRO:HD2	2.19	0.41
1:B:64:LYS:HE2	1:B:190:TYR:CE2	2.55	0.41
1:D:133:GLY:HA2	1:D:164:VAL:O	2.20	0.41
1:A:214:PHE:CD2	2:A:801:HEM:HAC	2.55	0.41
1:A:662:VAL:O	1:A:694:ILE:HA	2.20	0.41
1:C:438:CYS:HB2	1:C:439:PRO:HD2	2.03	0.41
1:A:81:ARG:HG2	1:A:87:LYS:HA	2.02	0.41
1:D:60:ASN:ND2	1:D:376:ASN:O	2.45	0.41
1:B:661:ALA:HB1	1:B:729:PHE:CZ	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:150:ALA:O	1:B:154:GLN:HG2	2.21	0.41
1:B:408:LEU:O	1:B:412:LEU:HG	2.21	0.41
1:B:750:GLY:HA2	1:D:682:HIS:CD2	2.56	0.41
1:C:393:PRO:HD2	1:C:415:TYR:CD1	2.55	0.41
3:C:925:HOH:O	1:D:480:THR:HG22	2.20	0.41
1:D:239:PHE:O	1:D:243:ILE:HG12	2.20	0.41
1:D:268:MET:HA	1:D:298:GLY:O	2.20	0.41
1:C:59:LYS:HB3	1:C:59:LYS:HE3	1.88	0.40
1:C:364:LEU:HD11	1:C:577:ASN:HB2	2.03	0.40
1:C:711:LYS:HA	1:C:711:LYS:HD3	1.87	0.40
1:D:159:LYS:O	1:D:161:PRO:HD3	2.21	0.40
1:D:411:ARG:HD2	2:D:801:HEM:CHC	2.51	0.40
1:A:275:HIS:CG	1:A:408:LEU:HB2	2.57	0.40
1:D:296:VAL:HB	1:D:333:GLU:HB2	2.03	0.40
1:C:128:HIS:CE1	1:C:169:VAL:HG22	2.56	0.40
1:C:214:PHE:CD2	2:C:801:HEM:CAC	3.04	0.40
1:B:173:ARG:NH2	1:D:228:GLU:OE1	2.52	0.40
1:D:519:HIS:NE2	1:D:742:GLU:OE2	2.39	0.40
1:B:641:GLN:H	1:B:641:GLN:CD	2.24	0.40
1:C:541:VAL:HG12	1:C:563:LEU:HD21	2.03	0.40
1:D:169:VAL:HG21	1:D:182:ILE:HB	2.04	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	715/750 (95%)	698 (98%)	17 (2%)	0	100	100
1	B	717/750 (96%)	697 (97%)	20 (3%)	0	100	100
1	C	719/750 (96%)	696 (97%)	23 (3%)	0	100	100
1	D	716/750 (96%)	697 (97%)	19 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	2867/3000 (96%)	2788 (97%)	79 (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	625/652 (96%)	619 (99%)	6 (1%)	76	82
1	B	626/652 (96%)	619 (99%)	7 (1%)	73	79
1	C	628/652 (96%)	620 (99%)	8 (1%)	69	75
1	D	625/652 (96%)	620 (99%)	5 (1%)	81	86
All	All	2504/2608 (96%)	2478 (99%)	26 (1%)	76	82

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	76	GLN
1	A	191	THR
1	A	237	ASP
1	A	286	CYS
1	A	583	ARG
1	A	683	TYR
1	B	191	THR
1	B	237	ASP
1	B	414	SER
1	B	617	MET
1	B	656	SER
1	B	683	TYR
1	B	724	ASP
1	C	151	GLU
1	C	237	ASP
1	C	283	GLU
1	C	416	ILE

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Mol	Chain	Res	Type
1	C	454	SER
1	C	628	LYS
1	C	683	TYR
1	C	724	ASP
1	D	59	LYS
1	D	237	ASP
1	D	265	SER
1	D	392	CYS
1	D	414	SER

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

Mol	Chain	Res	Type
1	D	682	HIS

### 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 monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

4 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
2	HEM	B	801	1	27,50,50	2.98	9 (33%)	17,82,82	1.45	3 (17%)
2	HEM	A	801	1	27,50,50	2.62	6 (22%)	17,82,82	1.41	2 (11%)
2	HEM	D	801	1	27,50,50	2.59	7 (25%)	17,82,82	1.31	2 (11%)
2	HEM	C	801	1	27,50,50	2.07	6 (22%)	17,82,82	1.30	3 (17%)

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	B	801	1	-	0/6/54/54	-
2	HEM	A	801	1	-	0/6/54/54	-
2	HEM	D	801	1	-	0/6/54/54	-
2	HEM	C	801	1	-	0/6/54/54	-

All (28) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	801	HEM	C1D-ND	-8.63	1.18	1.36
2	B	801	HEM	C4B-NB	-8.12	1.19	1.36
2	D	801	HEM	C4B-NB	-6.85	1.22	1.36
2	B	801	HEM	C1D-ND	-6.34	1.23	1.36
2	B	801	HEM	C1A-NA	5.97	1.48	1.36
2	C	801	HEM	C3B-C2B	-5.94	1.32	1.40
2	D	801	HEM	C1A-NA	5.37	1.47	1.36
2	B	801	HEM	C3B-C2B	-5.34	1.33	1.40
2	D	801	HEM	C1D-ND	-5.32	1.25	1.36
2	A	801	HEM	C4B-NB	-5.06	1.25	1.36
2	A	801	HEM	C3B-C2B	-4.48	1.34	1.40
2	A	801	HEM	C3C-C2C	-4.19	1.34	1.40
2	D	801	HEM	C4A-NA	-3.93	1.28	1.36
2	C	801	HEM	C4B-NB	-3.87	1.28	1.36
2	B	801	HEM	C3C-C2C	-3.80	1.35	1.40
2	A	801	HEM	C1A-NA	3.71	1.43	1.36
2	D	801	HEM	C3B-C2B	-3.68	1.35	1.40
2	D	801	HEM	C3C-C2C	-3.56	1.35	1.40
2	B	801	HEM	C4A-NA	-3.29	1.29	1.36
2	C	801	HEM	C3C-C2C	-3.01	1.36	1.40
2	C	801	HEM	C1A-NA	2.98	1.42	1.36
2	C	801	HEM	C1D-ND	-2.55	1.30	1.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	801	HEM	C3D-C2D	-2.37	1.30	1.37
2	B	801	HEM	C3B-CAB	-2.28	1.43	1.47
2	B	801	HEM	C3D-C2D	-2.21	1.31	1.37
2	D	801	HEM	C1C-C2C	-2.09	1.37	1.42
2	A	801	HEM	CAD-C3D	-2.08	1.48	1.52
2	B	801	HEM	CAD-C3D	-2.06	1.48	1.52

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	801	HEM	CBD-CAD-C3D	-3.57	105.90	112.48
2	A	801	HEM	CBD-CAD-C3D	-3.56	105.91	112.48
2	B	801	HEM	C3C-C4C-NC	-2.94	105.40	110.94
2	C	801	HEM	C3C-C4C-NC	-2.87	105.52	110.94
2	C	801	HEM	CBD-CAD-C3D	-2.80	107.32	112.48
2	B	801	HEM	C4A-C3A-C2A	2.65	108.84	107.00
2	D	801	HEM	C3C-C4C-NC	-2.63	105.97	110.94
2	D	801	HEM	CBD-CAD-C3D	-2.58	107.73	112.48
2	C	801	HEM	C3B-C4B-NB	-2.12	106.47	109.21
2	A	801	HEM	C3C-C4C-NC	-2.05	107.08	110.94

There are no chirality outliers.

There are no torsion outliers.

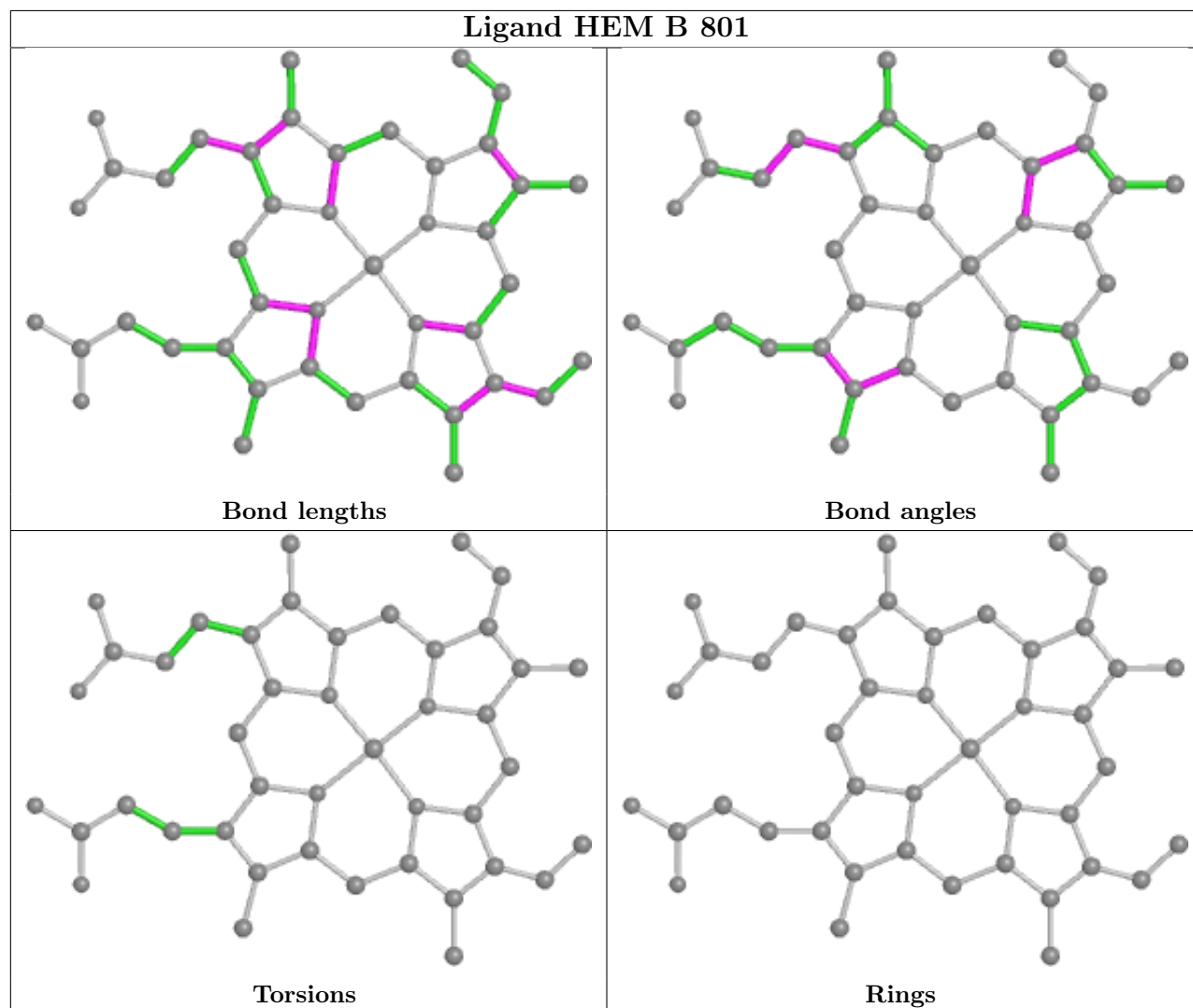
There are no ring outliers.

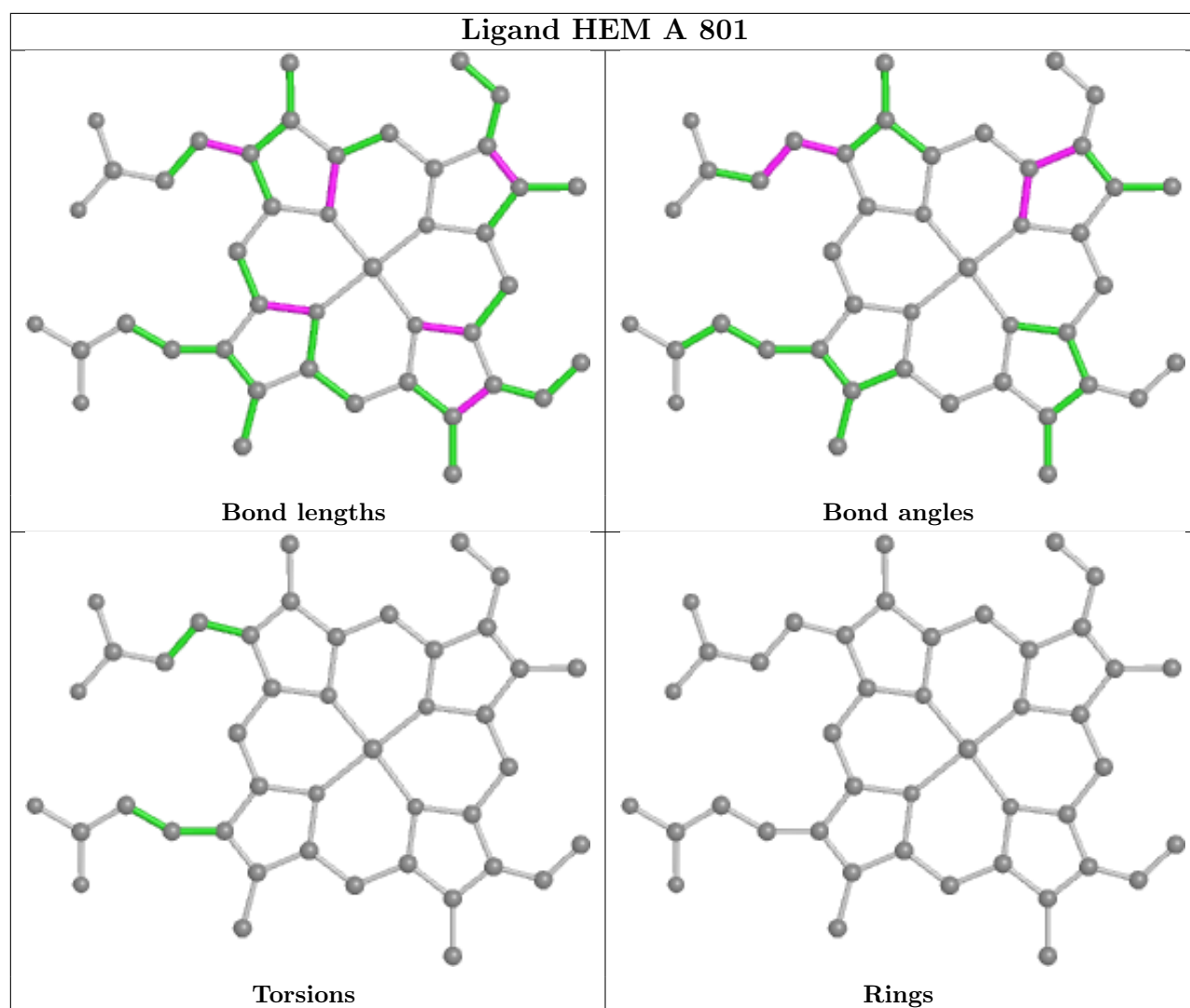
4 monomers are involved in 7 short contacts:

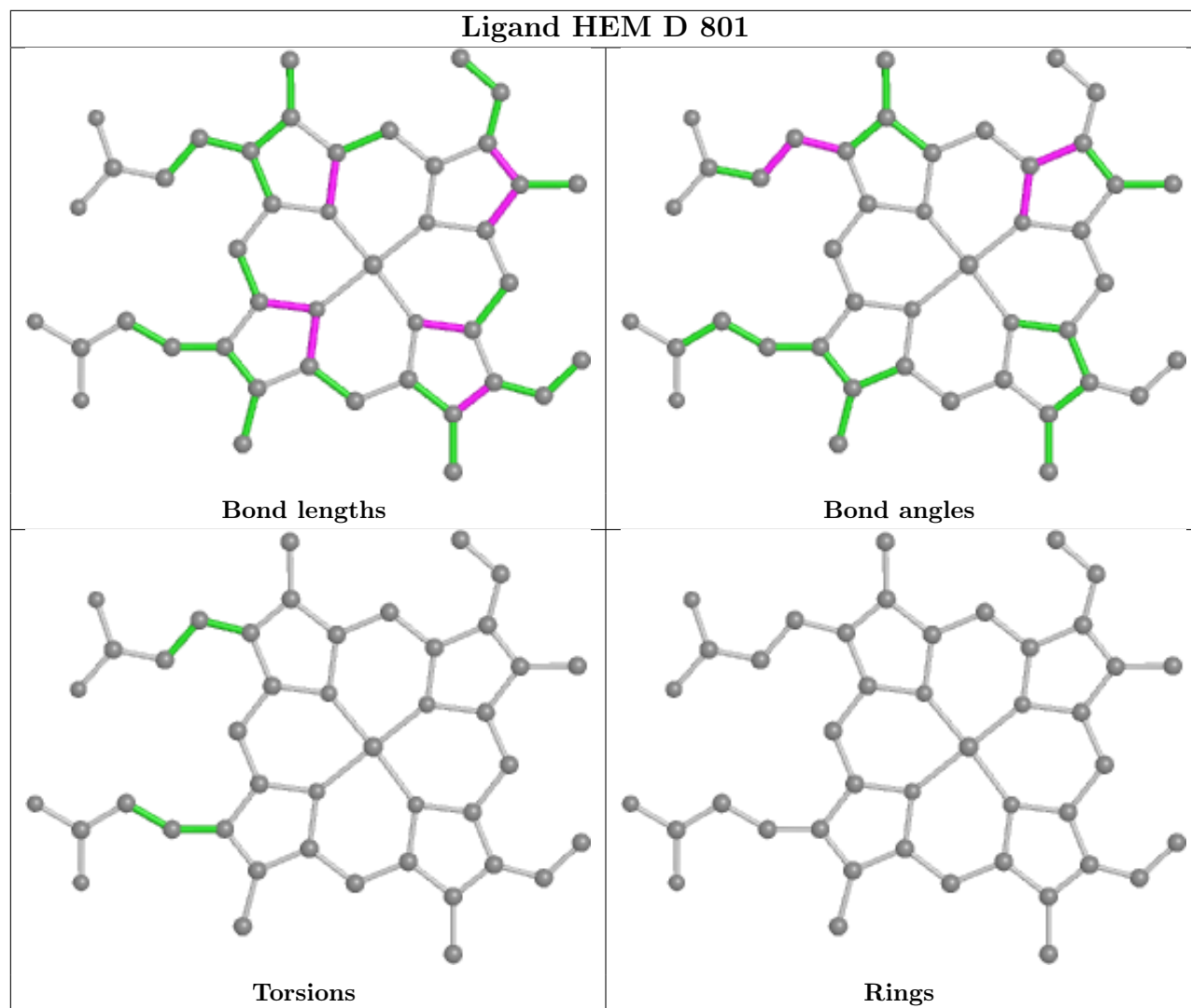
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	801	HEM	3	0
2	A	801	HEM	1	0
2	D	801	HEM	2	0
2	C	801	HEM	1	0

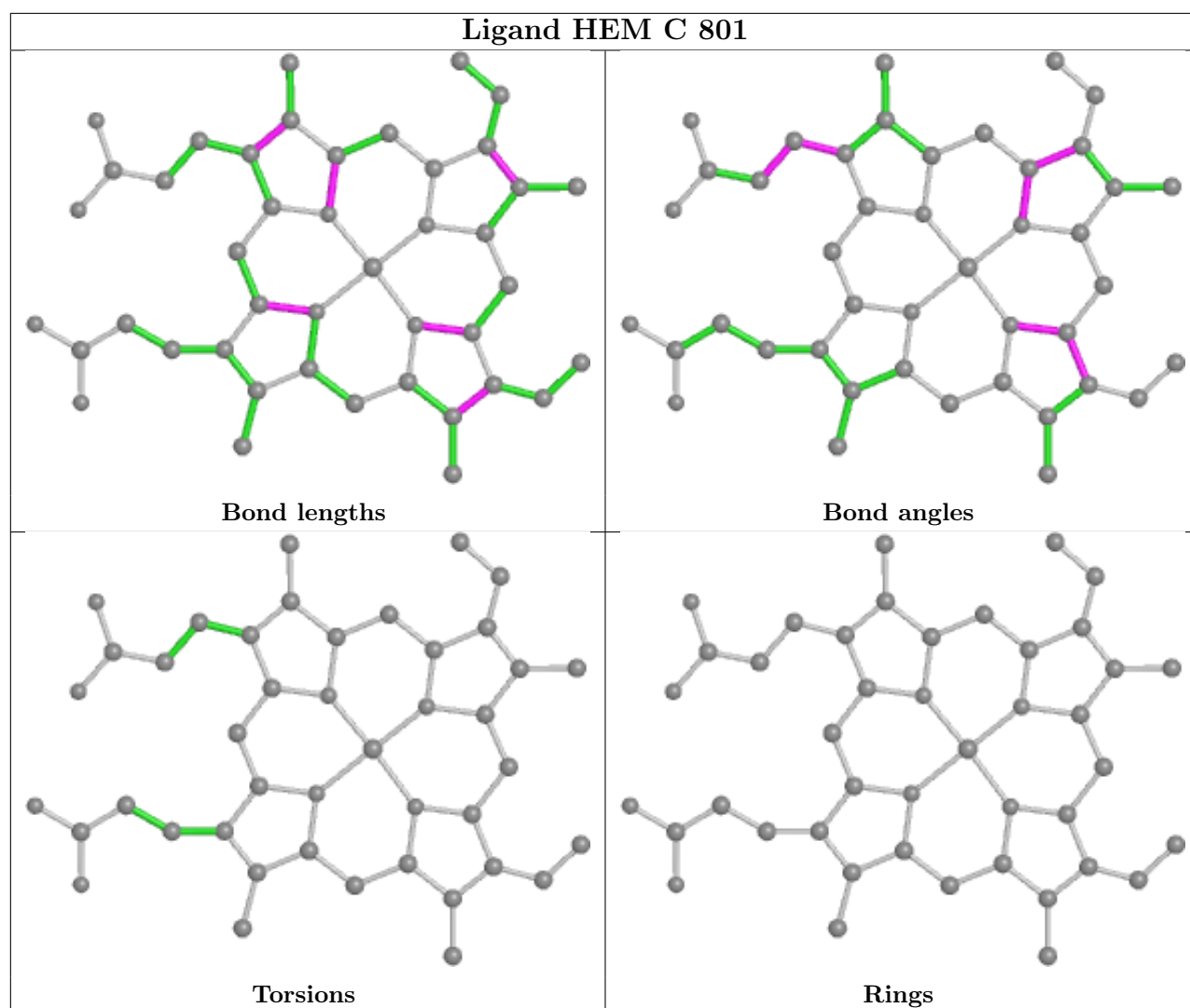
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

### 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	719/750 (95%)	-0.09	10 (1%) 75 78	20, 34, 50, 71	0
1	B	721/750 (96%)	-0.09	10 (1%) 75 78	19, 33, 51, 65	0
1	C	723/750 (96%)	0.01	18 (2%) 57 62	21, 35, 57, 67	0
1	D	720/750 (96%)	0.05	26 (3%) 42 49	21, 35, 65, 91	0
All	All	2883/3000 (96%)	-0.03	64 (2%) 62 66	19, 34, 55, 91	0

All (64) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	750	GLY	6.1
1	A	47	VAL	5.5
1	A	16	ASN	5.2
1	C	34	LYS	4.5
1	B	47	VAL	4.4
1	C	47	VAL	4.3
1	D	605	ALA	4.2
1	D	676	LYS	4.0
1	D	32	ALA	4.0
1	D	674	LEU	4.0
1	B	32	ALA	3.9
1	C	672	ASP	3.8
1	D	702	LYS	3.8
1	D	608	VAL	3.6
1	C	701	PHE	3.5
1	C	667	GLY	3.4
1	C	711	LYS	3.1
1	D	701	PHE	2.9
1	D	565	GLN	2.8
1	D	668	GLU	2.8
1	D	667	GLY	2.8

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Mol	Chain	Res	Type	RSRZ
1	C	718	VAL	2.8
1	C	664	ILE	2.7
1	A	73	PHE	2.7
1	C	724	ASP	2.6
1	D	700	ALA	2.6
1	D	698	GLY	2.6
1	C	668	GLU	2.5
1	B	701	PHE	2.5
1	C	706	ALA	2.5
1	A	612	ALA	2.5
1	D	562	LYS	2.4
1	A	592	HIS	2.4
1	D	642	GLY	2.4
1	A	30	ALA	2.4
1	D	721	LYS	2.4
1	C	702	LYS	2.3
1	A	312	GLY	2.3
1	C	623	TYR	2.3
1	C	721	LYS	2.2
1	B	720	ASP	2.2
1	D	611	ASP	2.2
1	A	17	LYS	2.2
1	C	696	LEU	2.2
1	A	179	VAL	2.2
1	B	702	LYS	2.2
1	C	33	GLU	2.2
1	D	675	MET	2.1
1	D	174	GLY	2.1
1	D	583	ARG	2.1
1	B	415	TYR	2.1
1	D	643	ASN	2.1
1	D	711	LYS	2.1
1	B	127	VAL	2.1
1	A	18	ALA	2.1
1	D	17	LYS	2.1
1	B	638	THR	2.1
1	D	16	ASN	2.1
1	B	73	PHE	2.1
1	C	712	PRO	2.0
1	C	725	LEU	2.0
1	B	592	HIS	2.0
1	D	669	ASP	2.0

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Mol	Chain	Res	Type	RSRZ
1	D	604	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 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.

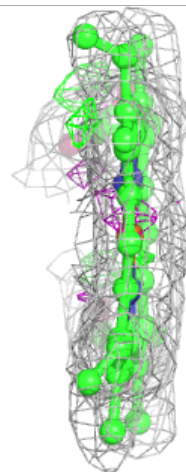
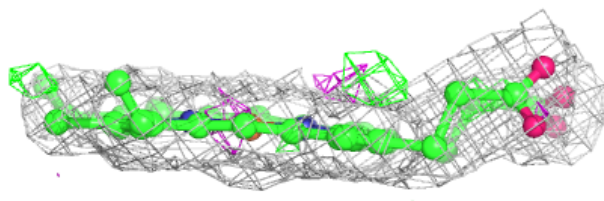
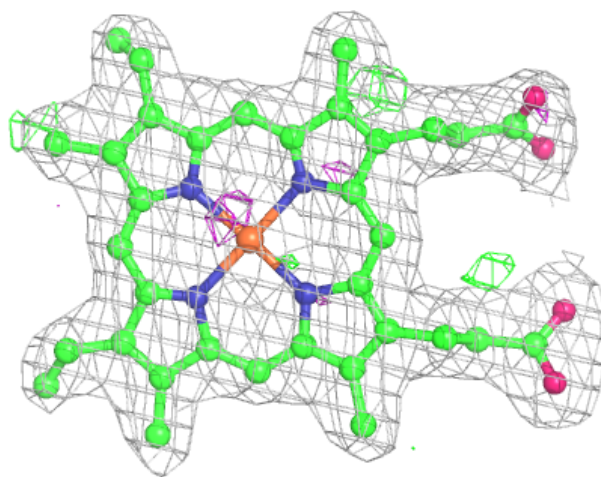
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
2	HEM	A	801	43/43	0.96	0.14	23,27,32,35	0
2	HEM	B	801	43/43	0.96	0.17	15,25,29,32	0
2	HEM	D	801	43/43	0.96	0.16	18,25,31,31	0
2	HEM	C	801	43/43	0.97	0.13	19,24,28,32	0

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.



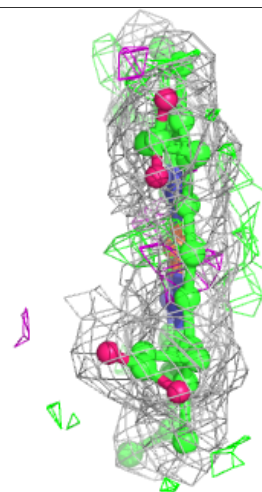
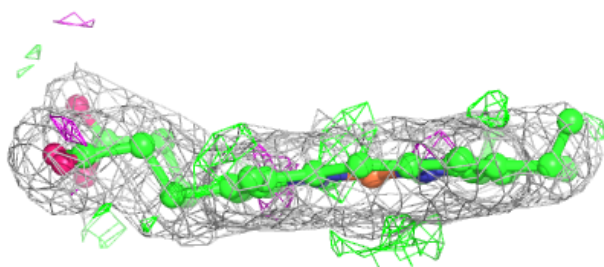
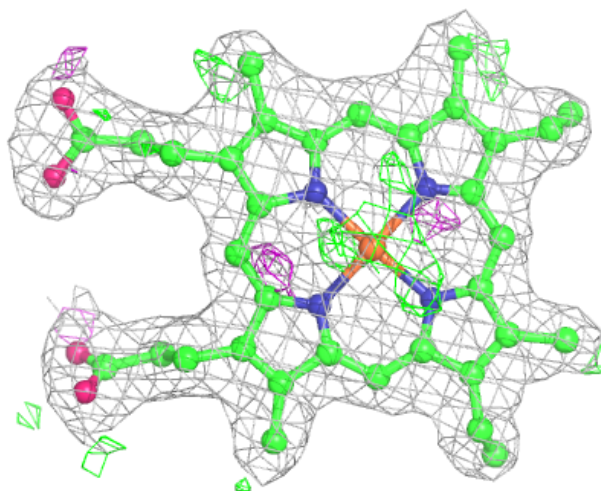
**Electron density around HEM A 801:**

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



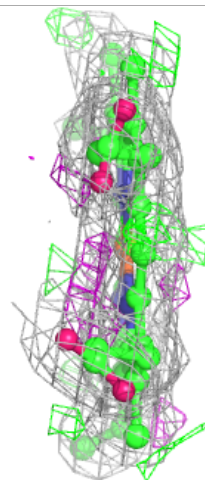
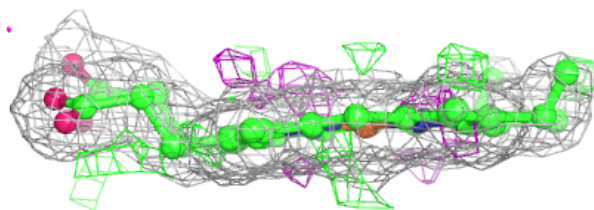
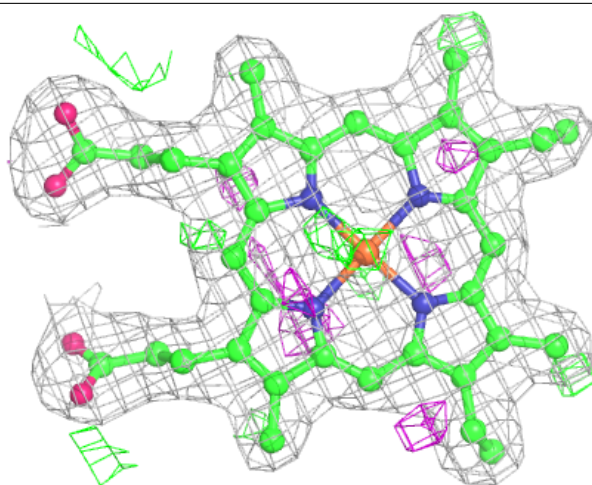
**Electron density around HEM B 801:**

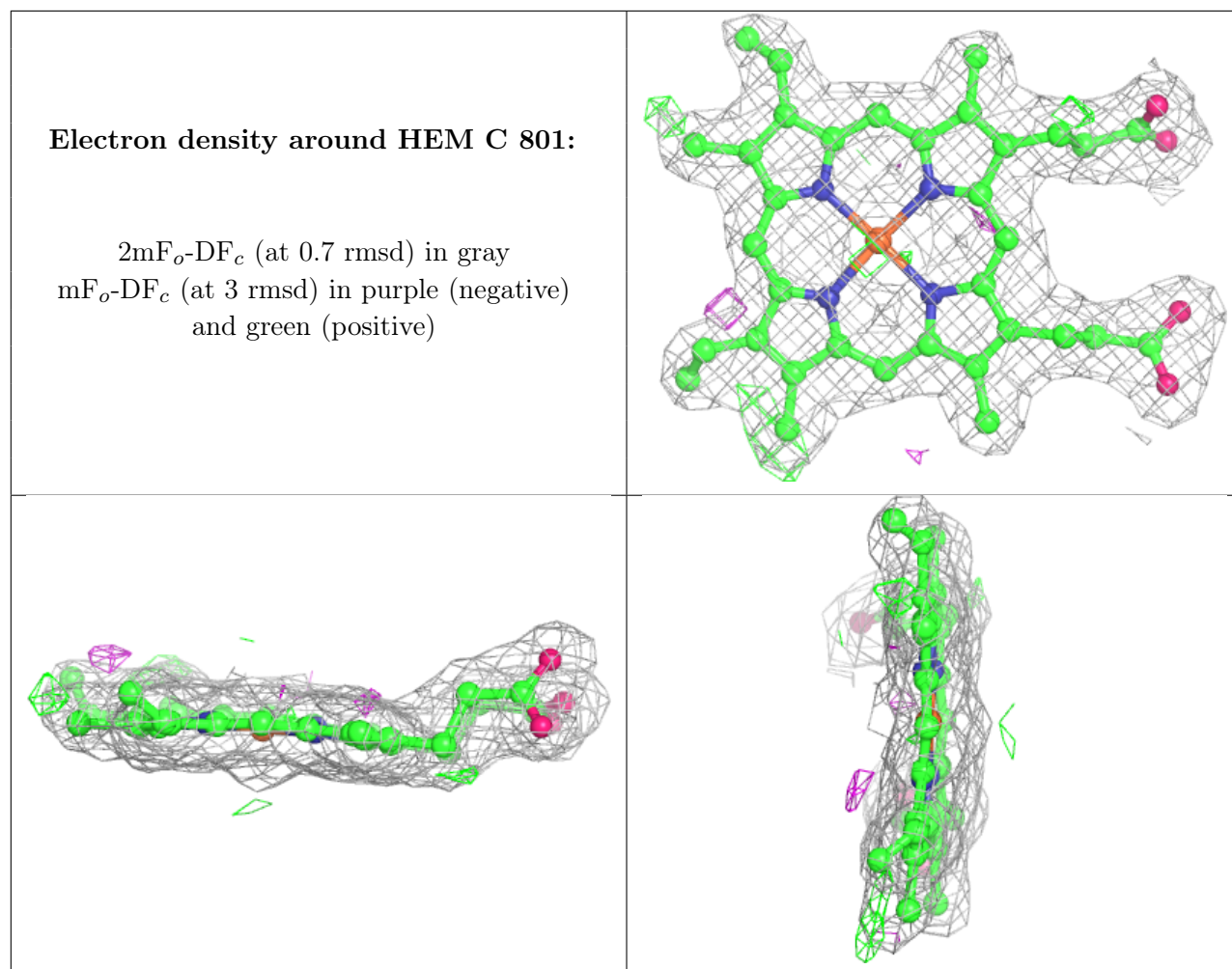
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



**Electron density around HEM 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 [i](#)

There are no such residues in this entry.