



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 11:45 AM GMT

PDB ID : 3PRC  
Title : PHOTOSYNTHETIC REACTION CENTER FROM RHODOPSEUDOMONAS VIRIDIS (QB-DEPLETED)  
Authors : Lancaster, C.R.D.; Michel, H.  
Deposited on : 1997-07-29  
Resolution : 2.40 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  
<http://wwpdb.org/validation/2016/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 : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

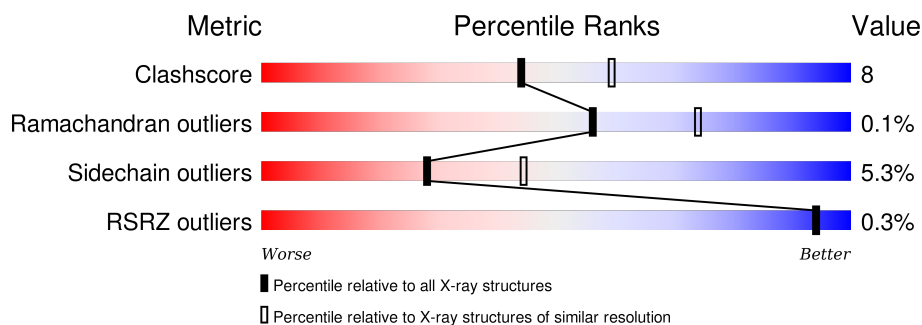
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.40 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	102246	3407 (2.40-2.40)
Ramachandran outliers	100387	3351 (2.40-2.40)
Sidechain outliers	100360	3352 (2.40-2.40)
RSRZ outliers	91569	2928 (2.40-2.40)

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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. 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	C	336	
2	L	273	
3	M	323	
4	H	258	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
11	NS5	M	600	-	-	-	X
12	LDA	L	702	-	-	-	X
12	LDA	L	707	-	-	-	X
12	LDA	M	704	-	-	-	X
12	LDA	M	705	-	-	-	X
12	LDA	M	706	-	-	-	X

## 2 Entry composition [i](#)

There are 13 unique types of molecules in this entry. The entry contains 10606 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 PHOTOSYNTHETIC REACTION CENTER.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	C	332	Total	C	N	O	S	38	1	0
			2607	1642	467	480	18			

- Molecule 2 is a protein called PHOTOSYNTHETIC REACTION CENTER.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	L	273	Total	C	N	O	S	14	2	0
			2193	1471	358	357	7			

- Molecule 3 is a protein called PHOTOSYNTHETIC REACTION CENTER.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	M	323	Total	C	N	O	S	14	1	0
			2566	1711	420	424	11			

- Molecule 4 is a protein called PHOTOSYNTHETIC REACTION CENTER.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	H	258	Total	C	N	O	S	129	1	0
			2028	1298	345	382	3			

- Molecule 5 is FE (II) ION (three-letter code: FE2) (formula: Fe).

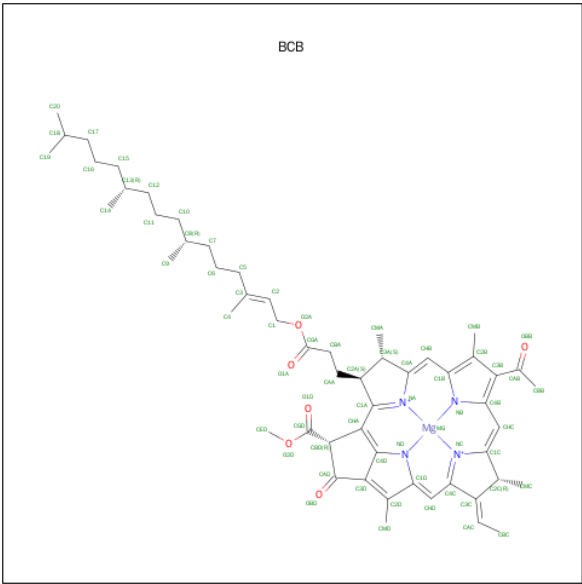
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	M	1	Total	Fe	0	0
			1	1		

- Molecule 6 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



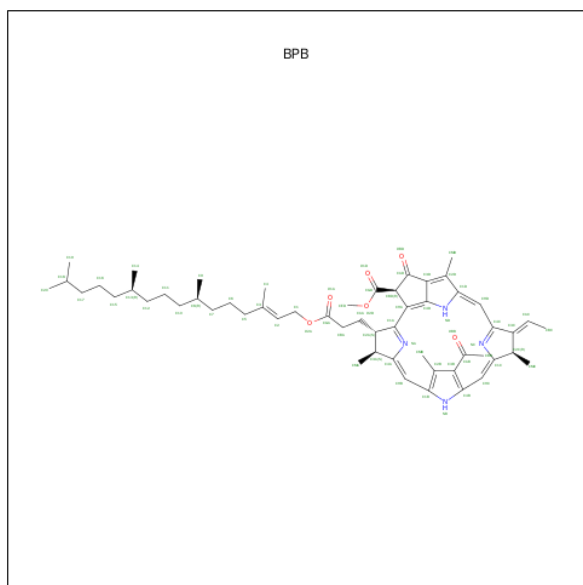
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	H	1	Total	O	S	0	0
			5	4	1		
6	M	1	Total	O	S	0	0
			5	4	1		
6	M	1	Total	O	S	0	0
			5	4	1		
6	M	1	Total	O	S	0	0
			5	4	1		

- Molecule 7 is BACTERIOCHLOROPHYLL B (three-letter code: BCB) (formula: C<sub>55</sub>H<sub>72</sub>MgN<sub>4</sub>O<sub>6</sub>).



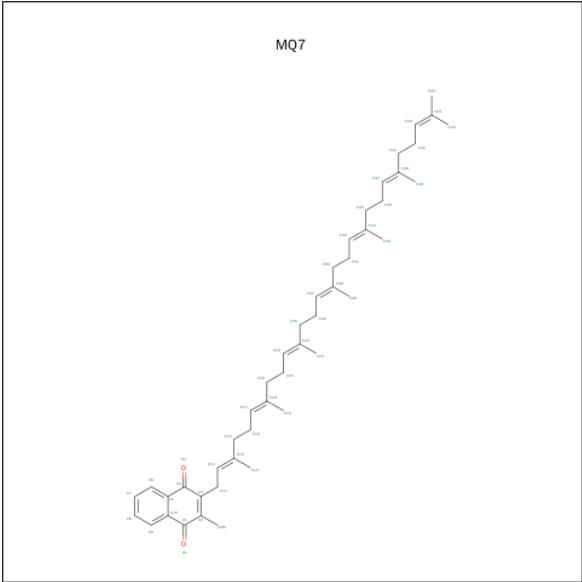
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
7	M	1	Total 66	C 55	Mg 1	N 4	O 6	0	0
7	L	1	Total 66	C 55	Mg 1	N 4	O 6	0	0
7	M	1	Total 66	C 55	Mg 1	N 4	O 6	0	0
7	L	1	Total 66	C 55	Mg 1	N 4	O 6	0	0

- Molecule 8 is BACTERIOPHEOPHYTIN B (three-letter code: BPB) (formula:  $C_{55}H_{74}N_4O_6$ ).



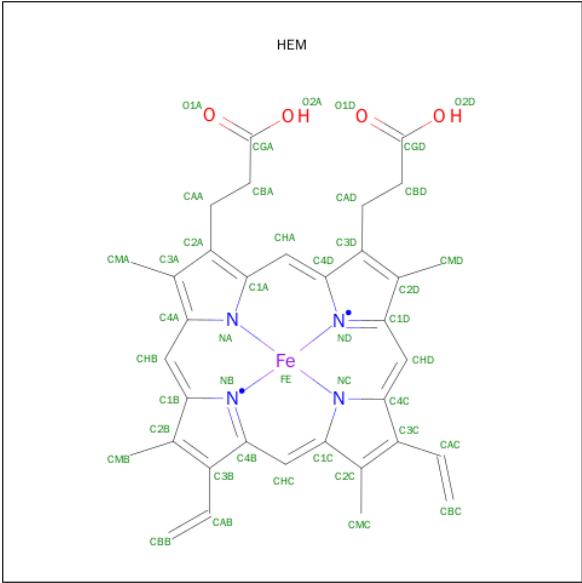
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
8	M	1	Total	C	N	O	7	0
			65	55	4	6		
8	L	1	Total	C	N	O	0	0
			65	55	4	6		

- Molecule 9 is MENAQUINONE-7 (three-letter code: MQ7) (formula:  $C_{46}H_{64}O_2$ ).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
9	M	1	Total	C	O		0	0
			48	46	2			

- Molecule 10 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (formula:  $C_{34}H_{32}FeN_4O_4$ ).



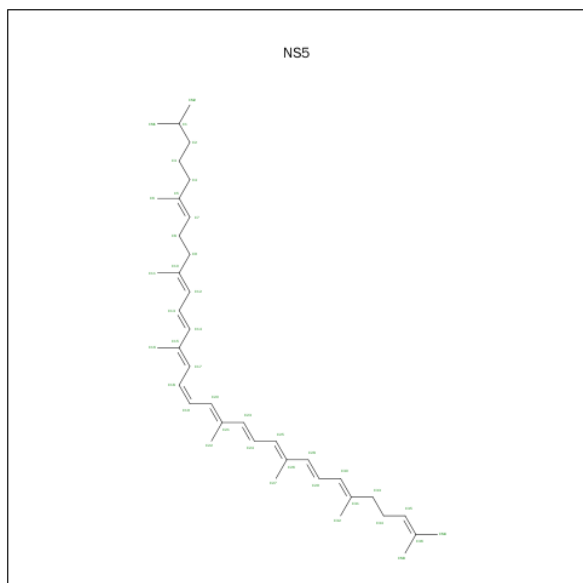
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
10	C	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
10	C	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		

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

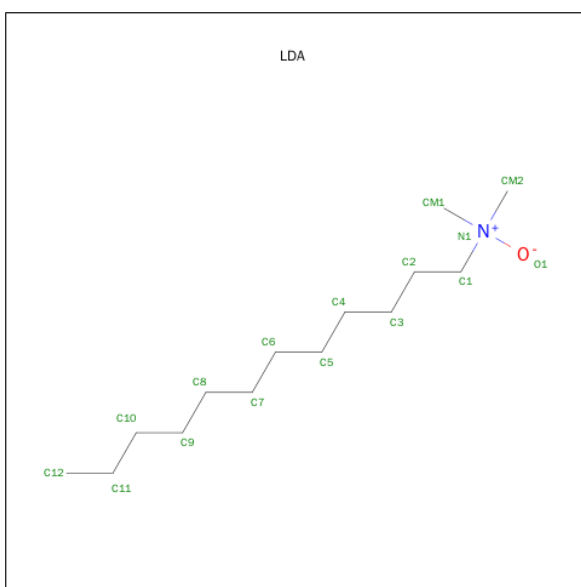
- Molecule 11 is 15-CIS-1,2-DIHYDRONEUROSPORENE (three-letter code: NS5) (formula:  $C_{40}H_{60}$ ).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
11	M	1	Total	C	9	0
			40	40		

- Molecule 12 is LAURYL DIMETHYLAMINE-N-OXIDE (three-letter code: LDA) (formula:  $C_{14}H_{31}NO$ ).





Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
12	H	1	Total	C	N	O	0	0
			16	14	1	1		
12	L	1	Total	C	N	O	0	0
			16	14	1	1		
12	H	1	Total	C	N	O	1	0
			16	14	1	1		
12	M	1	Total	C	N	O	3	0
			16	14	1	1		
12	M	1	Total	C	N	O	5	0
			16	14	1	1		
12	M	1	Total	C	N	O	4	0
			16	14	1	1		
12	L	1	Total	C	N	O	0	0
			16	14	1	1		

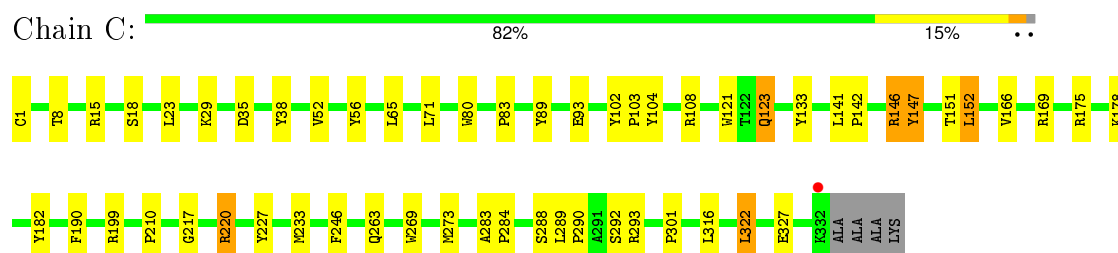
- Molecule 13 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
13	C	163	Total	O	0	0
			163	163		
13	H	100	Total	O	0	0
			100	100		
13	L	65	Total	O	0	0
			65	65		
13	M	97	Total	O	0	0
			97	97		

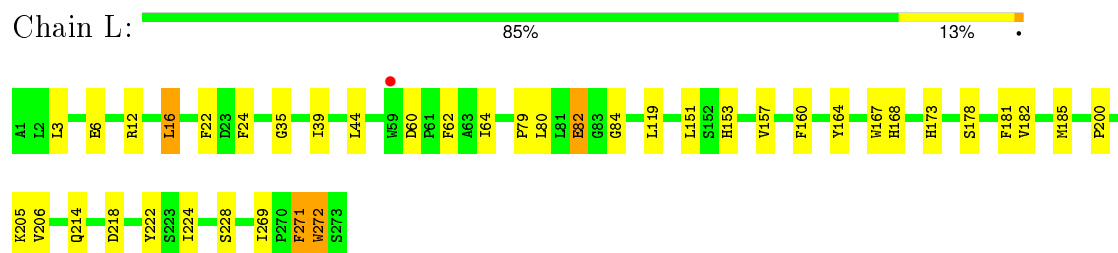
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

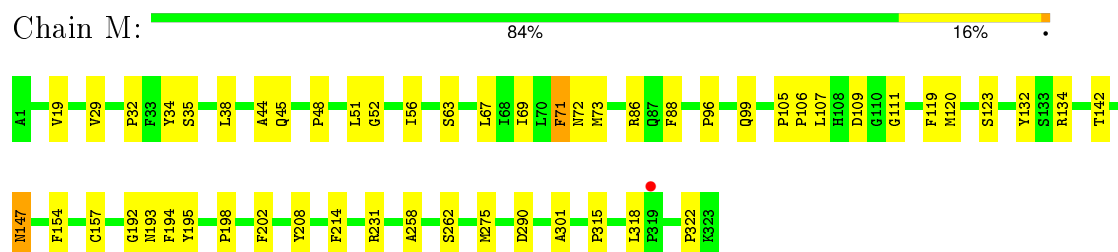
#### • Molecule 1: PHOTOSYNTHETIC REACTION CENTER



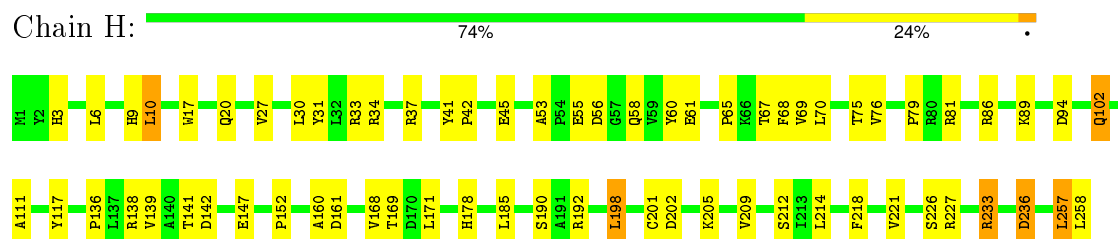
#### • Molecule 2: PHOTOSYNTHETIC REACTION CENTER



#### • Molecule 3: PHOTOSYNTHETIC REACTION CENTER



#### • Molecule 4: PHOTOSYNTHETIC REACTION CENTER



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	223.50 Å   223.50 Å   113.60 Å 90.00°   90.00°   90.00°	Depositor
Resolution (Å)	10.00 – 2.40 27.62 – 2.40	Depositor EDS
% Data completeness (in resolution range)	79.5 (10.00-2.40) 79.5 (27.62-2.40)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	0.09	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.69 (at 2.39 Å)	Xtriage
Refinement program	X-PLOR	Depositor
R, $R_{free}$	0.178   ,   0.215 0.166   ,   (Not available)	Depositor DCC
$R_{free}$ test set	No test flags present.	DCC
Wilson B-factor (Å <sup>2</sup> )	30.3	Xtriage
Anisotropy	0.030	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 89.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	0 of 89029 reflections	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	10606	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.46% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: LDA, BPB, BCB, FE2, SO4, MQ7, HEM, FME, NS5

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	C	0.51	0/2674	0.60	0/3645
2	L	0.53	0/2281	0.57	0/3112
3	M	0.52	0/2671	0.58	0/3653
4	H	0.53	0/2055	0.71	3/2807 (0.1%)
All	All	0.52	0/9681	0.61	3/13217 (0.0%)

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	5
2	L	0	1
3	M	0	1
All	All	0	7

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	H	53	ALA	CB-CA-C	-11.82	92.36	110.10
4	H	45	GLU	CB-CA-C	-7.85	94.70	110.40
4	H	53	ALA	N-CA-C	7.52	131.30	111.00

There are no chirality outliers.

All (7) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	147	TYR	Sidechain
1	C	182	TYR	Sidechain
1	C	190	PHE	Sidechain
1	C	227	TYR	Sidechain
1	C	89	TYR	Sidechain
2	L	164	TYR	Sidechain
3	M	119	PHE	Sidechain

## 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	C	2607	0	2575	33	0
2	L	2193	0	2122	31	0
3	M	2566	0	2460	37	0
4	H	2028	0	2029	38	0
5	M	1	0	0	0	0
6	H	5	0	0	0	0
6	M	15	0	0	0	0
7	L	132	0	144	9	0
7	M	132	0	144	13	0
8	L	65	0	74	5	0
8	M	65	0	74	12	0
9	M	48	0	64	0	0
10	C	172	0	120	1	0
11	M	40	0	60	0	0
12	H	32	0	62	0	0
12	L	32	0	62	10	0
12	M	48	0	93	2	0
13	C	163	0	0	5	0
13	H	100	0	0	2	0
13	L	65	0	0	2	0
13	M	97	0	0	0	0
All	All	10606	0	10083	157	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 8.

All (157) 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)
8:M:401:BPB:HHC	8:M:401:BPB:HBBB	1.50	0.89
8:L:402:BPB:HHC	8:L:402:BPB:HBBB	1.53	0.89
7:M:805:BCB:HHC	7:M:805:BCB:HBB2	1.58	0.82
1:C:121:TRP:HA	1:C:123[A]:GLN:HE21	1.43	0.81
8:L:402:BPB:HHC	8:L:402:BPB:CBB	2.10	0.81
3:M:29:VAL:HG21	3:M:51:LEU:HD22	1.64	0.79
1:C:152:LEU:HD22	1:C:175:ARG:HA	1.66	0.75
2:L:79:PRO:HG2	2:L:82:GLU:HG3	1.72	0.71
7:M:806:BCB:HAA2	7:M:806:BCB:HBD	1.74	0.69
1:C:123[A]:GLN:HG3	1:C:269:TRP:CE3	2.28	0.69
2:L:178:SER:O	2:L:182:VAL:HG23	1.95	0.67
2:L:181:PHE:HB3	8:M:401:BPB:CBB	2.24	0.66
4:H:55:GLU:HB2	4:H:58:GLN:HG3	1.78	0.65
1:C:217:GLY:O	1:C:220:ARG:HG3	1.97	0.65
1:C:102:TYR:CD2	1:C:103:PRO:HD3	2.33	0.64
12:L:707:LDA:HM23	13:L:761:HOH:O	1.99	0.63
3:M:71:PHE:HB3	12:M:706:LDA:H62	1.80	0.63
7:M:805:BCB:CBB	7:M:805:BCB:HHC	2.30	0.62
1:C:290:PRO:HG2	1:C:293:ARG:HG2	1.81	0.62
4:H:160:ALA:HB3	4:H:214:LEU:HD23	1.80	0.62
1:C:152:LEU:HD21	1:C:178:LYS:HG3	1.82	0.62
1:C:121:TRP:HA	1:C:123[A]:GLN:NE2	2.13	0.62
4:H:161:ASP:HB3	4:H:214:LEU:HD22	1.81	0.62
12:L:702:LDA:HM23	13:L:772:HOH:O	1.98	0.61
2:L:181:PHE:HB3	8:M:401:BPB:HBBA	1.81	0.61
3:M:315:PRO:HA	3:M:318:LEU:HG	1.82	0.60
2:L:62:PHE:HE2	12:L:702:LDA:HM21	1.66	0.59
4:H:117:TYR:HB2	4:H:236:ASP:HB3	1.85	0.59
1:C:65:LEU:HD11	1:C:327:GLU:HG2	1.85	0.59
3:M:32:PRO:HG3	3:M:48:PRO:HD3	1.83	0.59
2:L:218:ASP:HB3	3:M:134:ARG:HD2	1.85	0.59
3:M:63:SER:O	3:M:67:LEU:HG	2.03	0.58
3:M:29:VAL:CG2	3:M:51:LEU:HD22	2.34	0.58
3:M:231:ARG:HD2	13:H:812:HOH:O	2.03	0.57
8:M:401:BPB:CHC	8:M:401:BPB:HBBB	2.28	0.57
2:L:62:PHE:CE2	12:L:702:LDA:HM21	2.40	0.57
2:L:12:ARG:HD3	4:H:102:GLN:NE2	2.19	0.57
2:L:181:PHE:CD2	8:M:401:BPB:HBB	2.40	0.56
7:L:304:BCB:HMB1	7:L:304:BCB:HBB2	1.87	0.56
3:M:202:PHE:CE2	4:H:20:GLN:HG2	2.40	0.56
7:L:304:BCB:HMB2	8:L:402:BPB:HMBA	1.88	0.56
2:L:35:GLY:O	2:L:39:ILE:HG12	2.05	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
12:L:702:LDA:H32	13:H:894:HOH:O	2.06	0.56
3:M:96:PRO:HG3	3:M:105:PRO:HB3	1.87	0.56
2:L:153:HIS:O	2:L:157:VAL:HG23	2.06	0.55
1:C:80:TRP:CD1	1:C:133:TYR:HB2	2.42	0.55
2:L:80:LEU:HA	2:L:84:GLY:HA3	1.88	0.55
7:M:805:BCB:H61	8:M:401:BPB:HMA	1.89	0.54
4:H:136:PRO:HG2	4:H:139:VAL:HG23	1.90	0.54
4:H:6:LEU:HB2	4:H:10:LEU:HB3	1.89	0.54
7:M:806:BCB:H203	8:M:401:BPB:H4B	1.91	0.53
2:L:185:MET:SD	7:M:805:BCB:H41	2.47	0.53
7:M:806:BCB:HAA2	7:M:806:BCB:CBD	2.34	0.53
2:L:200:PRO:HG3	2:L:205:LYS:O	2.08	0.53
1:C:210:PRO:HB2	4:H:3:HIS:HD2	1.74	0.53
4:H:56:ASP:HB3	4:H:60:TYR:CE2	2.44	0.53
3:M:275:MET:HG2	8:M:401:BPB:HBCA	1.90	0.53
7:L:302:BCB:HBB3	7:L:302:BCB:HMB1	1.89	0.53
4:H:86:ARG:NH2	4:H:111:ALA:O	2.37	0.53
4:H:136:PRO:HG2	4:H:138:ARG:HG2	1.92	0.52
4:H:233:ARG:HG3	4:H:233:ARG:O	2.09	0.52
12:L:702:LDA:H71	3:M:301:ALA:CB	2.40	0.52
1:C:147:TYR:OH	1:C:301:PRO:HG3	2.10	0.52
7:L:302:BCB:OBB	7:L:302:BCB:HHC	2.09	0.51
4:H:190:SER:HB3	4:H:192:ARG:HG2	1.90	0.51
2:L:214:GLN:NE2	3:M:19:VAL:H	2.08	0.51
4:H:86:ARG:NH2	4:H:111:ALA:HB3	2.25	0.51
7:L:304:BCB:HMB1	7:L:304:BCB:CBB	2.41	0.50
1:C:289:LEU:HD22	1:C:293:ARG:HG3	1.93	0.50
4:H:152:PRO:O	4:H:168:VAL:HB	2.12	0.50
2:L:182:VAL:HG22	7:M:805:BCB:H12	1.94	0.50
1:C:93:GLU:HB2	13:C:470:HOH:O	2.10	0.50
7:L:302:BCB:HMB1	7:L:302:BCB:CBB	2.42	0.49
3:M:99:GLN:OE1	3:M:99:GLN:HA	2.11	0.49
2:L:224:ILE:HG12	2:L:228:SER:HB2	1.93	0.49
1:C:121:TRP:CG	1:C:273:MET:HG3	2.48	0.49
1:C:210:PRO:HB2	4:H:3:HIS:CD2	2.48	0.49
3:M:147:ASN:HD22	8:M:401:BPB:HMDA	1.77	0.49
2:L:206:VAL:HG21	4:H:68:PHE:HB3	1.93	0.49
4:H:65:PRO:HA	4:H:79:PRO:HD2	1.94	0.48
4:H:218:PHE:HA	4:H:221:VAL:HG23	1.95	0.48
3:M:34:TYR:HA	3:M:44:ALA:O	2.14	0.48
1:C:102:TYR:CG	1:C:103:PRO:HD3	2.48	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:M:32:PRO:HB3	3:M:45:GLN:HG3	1.95	0.48
1:C:233:MET:HB3	10:C:339:HEM:C4B	2.49	0.48
3:M:52:GLY:O	3:M:56:ILE:HD12	2.14	0.47
2:L:168:HIS:CE1	7:L:302:BCB:HMC2	2.50	0.47
8:M:401:BPB:HHC	8:M:401:BPB:CBB	2.31	0.47
3:M:29:VAL:HG21	3:M:51:LEU:CD2	2.40	0.47
1:C:220:ARG:NH2	13:C:359:HOH:O	2.47	0.47
2:L:269:ILE:HG22	2:L:271:PHE:HD1	1.79	0.47
1:C:146:ARG:HA	1:C:146:ARG:HD2	1.64	0.47
12:L:702:LDA:H71	3:M:301:ALA:HB2	1.96	0.47
7:M:806:BCB:OBB	7:M:806:BCB:HHC	2.16	0.46
1:C:83:PRO:HD2	13:C:367:HOH:O	2.15	0.46
3:M:132:TYR:CE1	3:M:142:THR:HG21	2.51	0.46
2:L:272:TRP:CE2	3:M:86:ARG:HG3	2.51	0.46
1:C:283:ALA:HB3	1:C:284:PRO:HD3	1.98	0.45
2:L:269:ILE:CG2	2:L:271:PHE:HD1	2.30	0.45
4:H:138:ARG:HG3	4:H:139:VAL:HG23	1.99	0.45
4:H:67:THR:HA	4:H:76:VAL:O	2.16	0.45
3:M:107:LEU:HA	3:M:111:GLY:HA3	1.98	0.44
4:H:37:ARG:HG2	4:H:41:TYR:CE1	2.52	0.44
4:H:34:ARG:HG2	4:H:61:GLU:O	2.17	0.44
1:C:283:ALA:N	1:C:284:PRO:CD	2.81	0.44
2:L:60:ASP:O	2:L:64:ILE:HG13	2.16	0.44
7:L:304:BCB:HHC	7:L:304:BCB:OBB	2.16	0.44
1:C:141:LEU:HD12	1:C:142:PRO:HD2	1.98	0.44
1:C:322:LEU:HA	13:C:469:HOH:O	2.17	0.44
8:L:402:BPB:NC	8:L:402:BPB:ND	2.66	0.44
2:L:151:LEU:HD21	12:L:702:LDA:H111	2.00	0.44
1:C:52:VAL:HB	1:C:56:TYR:CD2	2.53	0.44
1:C:104:TYR:HE1	1:C:108:ARG:NH2	2.16	0.44
4:H:33:ARG:HA	4:H:33:ARG:HD2	1.82	0.44
3:M:35:SER:HB3	3:M:38:LEU:HB3	2.00	0.43
3:M:72:ASN:OD1	12:M:706:LDA:H12	2.18	0.43
3:M:120:MET:HG3	7:M:806:BCB:H172	2.00	0.43
4:H:67:THR:CG2	4:H:75:THR:HB	2.47	0.43
2:L:3:LEU:HB2	2:L:6:GLU:HB2	1.99	0.43
1:C:246:PHE:CZ	1:C:263:GLN:HG2	2.54	0.43
3:M:195:TYR:CZ	7:M:806:BCB:HMC2	2.53	0.43
2:L:222:TYR:O	12:L:707:LDA:HM13	2.19	0.43
3:M:67:LEU:O	3:M:71:PHE:HB2	2.18	0.43
2:L:167:TRP:HE1	2:L:173:HIS:CD2	2.37	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:L:206:VAL:CG2	4:H:68:PHE:HB3	2.48	0.43
3:M:258:ALA:HB1	3:M:262:SER:OG	2.17	0.43
7:M:805:BCB:H203	7:M:805:BCB:H162	1.92	0.43
8:L:402:BPB:HBBA	3:M:208:TYR:HB3	2.00	0.43
8:M:401:BPB:H4	8:M:401:BPB:H6	1.77	0.43
1:C:8:THR:HB	1:C:23:LEU:HB2	2.01	0.43
4:H:27:VAL:O	4:H:31:TYR:HB3	2.19	0.42
1:C:52:VAL:HB	1:C:56:TYR:HD2	1.83	0.42
4:H:86:ARG:HH21	4:H:111:ALA:HB3	1.84	0.42
4:H:138:ARG:CG	4:H:139:VAL:HG23	2.49	0.42
4:H:70:LEU:HD11	4:H:76:VAL:HG23	2.00	0.42
2:L:185:MET:SD	7:M:805:BCB:C4	3.07	0.42
3:M:69:ILE:HG22	3:M:73:MET:SD	2.59	0.42
1:C:71:LEU:HD23	1:C:71:LEU:HA	1.88	0.42
8:M:401:BPB:CHC	8:M:401:BPB:CBB	2.96	0.42
4:H:257:LEU:HA	4:H:257:LEU:HD22	1.80	0.42
7:L:302:BCB:H41	7:L:302:BCB:H62	1.90	0.41
4:H:37:ARG:HG2	4:H:41:TYR:CZ	2.56	0.41
2:L:16:LEU:HA	2:L:16:LEU:HD12	1.71	0.41
3:M:192:GLY:O	3:M:193:ASN:HB3	2.20	0.41
4:H:202:ASP:HB3	4:H:209:VAL:HB	2.01	0.41
1:C:146:ARG:NH2	13:C:344:HOH:O	2.50	0.41
3:M:73:MET:HE1	3:M:88:PHE:CE1	2.56	0.41
1:C:35:ASP:HB3	1:C:316:LEU:HA	2.03	0.41
3:M:198:PRO:HB3	4:H:17:TRP:CZ3	2.56	0.41
4:H:152:PRO:HD2	4:H:171:LEU:HD11	2.02	0.41
4:H:142:ASP:N	4:H:142:ASP:OD1	2.54	0.41
12:L:702:LDA:H62	12:L:702:LDA:H11	2.02	0.40
4:H:198:LEU:O	4:H:201:CYS:HB2	2.21	0.40
2:L:22:PHE:HA	2:L:24:PHE:CE1	2.55	0.40
3:M:106:PRO:HG2	3:M:109:ASP:HB3	2.04	0.40
3:M:120:MET:O	3:M:123:SER:HB3	2.21	0.40
3:M:154:PHE:O	3:M:157:CYS:HB2	2.22	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles

### 5.3.1 Protein backbone

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	C	331/336 (98%)	320 (97%)	11 (3%)	0	100	100
2	L	273/273 (100%)	264 (97%)	9 (3%)	0	100	100
3	M	322/323 (100%)	308 (96%)	13 (4%)	1 (0%)	46	63
4	H	256/258 (99%)	247 (96%)	9 (4%)	0	100	100
All	All	1182/1190 (99%)	1139 (96%)	42 (4%)	1 (0%)	56	74

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	M	322	PRO

### 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	C	281/282 (100%)	264 (94%)	17 (6%)	24	37
2	L	220/218 (101%)	213 (97%)	7 (3%)	46	68
3	M	250/249 (100%)	245 (98%)	5 (2%)	63	81
4	H	212/212 (100%)	189 (89%)	23 (11%)	8	11
All	All	963/961 (100%)	911 (95%)	52 (5%)	28	43

All (52) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	1	CYS
1	C	15	ARG
1	C	18	SER
1	C	29	LYS
1	C	38	TYR
1	C	123[A]	GLN
1	C	123[B]	GLN
1	C	146	ARG
1	C	151	THR
1	C	152	LEU
1	C	166	VAL
1	C	169	ARG
1	C	199	ARG
1	C	220	ARG
1	C	288	SER
1	C	292	SER
1	C	322	LEU
2	L	16	LEU
2	L	44	LEU
2	L	82	GLU
2	L	119	LEU
2	L	160	PHE
2	L	271	PHE
2	L	272	TRP
3	M	71	PHE
3	M	147	ASN
3	M	194	PHE
3	M	214	PHE
3	M	290	ASP
4	H	9	HIS
4	H	10	LEU
4	H	30	LEU
4	H	42	PRO
4	H	69	VAL
4	H	81	ARG
4	H	89	LYS
4	H	94	ASP
4	H	102	GLN
4	H	141	THR
4	H	147	GLU
4	H	169	THR
4	H	178	HIS
4	H	185	LEU

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Mol	Chain	Res	Type
4	H	198	LEU
4	H	205	LYS
4	H	212	SER
4	H	226	SER
4	H	227	ARG
4	H	233	ARG
4	H	236	ASP
4	H	257	LEU
4	H	258	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	37	GLN
1	C	302	GLN
2	L	183	ASN
2	L	214	GLN
2	L	239	ASN
3	M	16	HIS
3	M	147	ASN
4	H	102	GLN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

2 non-standard protein/DNA/RNA residues 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	$\# Z  > 2$
4	FME	H	1[A]	4	8,9,10	0.61	0	6,9,11	3.28	3 (50%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
4	FME	H	1[B]	4	8,9,10	0.55	0	6,9,11	4.31	2 (33%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	FME	H	1[A]	4	-	0/6/9/11	0/0/0/0
4	FME	H	1[B]	4	-	0/6/9/11	0/0/0/0

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	H	1[B]	FME	O1-CN-N	-8.80	112.09	124.76
4	H	1[A]	FME	O1-CN-N	-6.61	115.25	124.76
4	H	1[B]	FME	CA-N-CN	-5.33	114.62	122.82
4	H	1[A]	FME	CA-N-CN	-3.81	116.96	122.82
4	H	1[A]	FME	O-C-CA	-2.13	119.82	125.44

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

Of 24 ligands modelled in this entry, 1 is monoatomic - leaving 23 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected

value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
10	HEM	C	337	1	30,50,50	2.73	9 (30%)	24,82,82	3.00	10 (41%)
10	HEM	C	338	1	30,50,50	2.58	11 (36%)	24,82,82	3.33	9 (37%)
10	HEM	C	339	1	30,50,50	2.66	9 (30%)	24,82,82	3.08	8 (33%)
10	HEM	C	340	1	30,50,50	2.70	10 (33%)	24,82,82	2.96	9 (37%)
12	LDA	H	701	-	15,15,15	4.90	3 (20%)	16,17,17	1.07	2 (12%)
12	LDA	H	703	-	15,15,15	4.39	3 (20%)	16,17,17	0.64	0
6	SO4	H	801	-	4,4,4	1.39	0	6,6,6	0.55	0
7	BCB	L	302	2	56,74,74	1.55	6 (10%)	57,115,115	1.81	10 (17%)
7	BCB	L	304	2	56,74,74	1.56	4 (7%)	57,115,115	2.03	10 (17%)
8	BPB	L	402	-	63,70,70	1.40	8 (12%)	63,101,101	1.81	9 (14%)
12	LDA	L	702	-	15,15,15	4.16	3 (20%)	16,17,17	0.79	0
12	LDA	L	707	-	15,15,15	3.61	2 (13%)	16,17,17	0.77	0
8	BPB	M	401	-	63,70,70	1.37	8 (12%)	63,101,101	1.85	9 (14%)
9	MQ7	M	501	-	49,49,49	1.70	11 (22%)	62,63,63	1.43	10 (16%)
11	NS5	M	600	-	39,39,39	0.73	0	44,46,46	1.12	5 (11%)
12	LDA	M	704	-	15,15,15	4.39	4 (26%)	16,17,17	0.62	0
12	LDA	M	705	-	15,15,15	4.11	4 (26%)	16,17,17	0.78	0
12	LDA	M	706	-	15,15,15	4.34	4 (26%)	16,17,17	0.62	0
6	SO4	M	802	-	4,4,4	2.08	2 (50%)	6,6,6	1.17	1 (16%)
6	SO4	M	803	-	4,4,4	1.32	0	6,6,6	0.56	0
6	SO4	M	804	-	4,4,4	1.98	1 (25%)	6,6,6	0.45	0
7	BCB	M	805	3	56,74,74	1.46	7 (12%)	57,115,115	1.98	11 (19%)
7	BCB	M	806	3	56,74,74	1.40	7 (12%)	57,115,115	2.25	13 (22%)

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
10	HEM	C	337	1	-	0/10/54/54	0/0/8/8
10	HEM	C	338	1	-	0/10/54/54	0/0/8/8
10	HEM	C	339	1	-	0/10/54/54	0/0/8/8
10	HEM	C	340	1	-	0/10/54/54	0/0/8/8
12	LDA	H	701	-	-	0/13/13/13	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
12	LDA	H	703	-	-	0/13/13/13	0/0/0/0
6	SO4	H	801	-	-	0/0/0/0	0/0/0/0
7	BCB	L	302	2	-	0/37/137/137	0/0/9/9
7	BCB	L	304	2	-	0/37/137/137	0/0/9/9
8	BPB	L	402	-	-	0/46/105/105	0/1/6/6
12	LDA	L	702	-	-	0/13/13/13	0/0/0/0
12	LDA	L	707	-	-	0/13/13/13	0/0/0/0
8	BPB	M	401	-	-	0/46/105/105	0/1/6/6
9	MQ7	M	501	-	-	0/41/61/61	0/2/2/2
11	NS5	M	600	-	-	0/43/43/43	0/0/0/0
12	LDA	M	704	-	-	0/13/13/13	0/0/0/0
12	LDA	M	705	-	-	0/13/13/13	0/0/0/0
12	LDA	M	706	-	-	0/13/13/13	0/0/0/0
6	SO4	M	802	-	-	0/0/0/0	0/0/0/0
6	SO4	M	803	-	-	0/0/0/0	0/0/0/0
6	SO4	M	804	-	-	0/0/0/0	0/0/0/0
7	BCB	M	805	3	-	0/37/137/137	0/0/9/9
7	BCB	M	806	3	-	0/37/137/137	0/0/9/9

All (116) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
12	H	701	LDA	O1-N1	-18.45	1.21	1.39
12	H	703	LDA	O1-N1	-16.61	1.23	1.39
12	M	704	LDA	O1-N1	-16.16	1.24	1.39
12	M	706	LDA	O1-N1	-16.11	1.24	1.39
12	L	702	LDA	O1-N1	-15.39	1.24	1.39
12	M	705	LDA	O1-N1	-15.14	1.25	1.39
12	L	707	LDA	O1-N1	-13.10	1.27	1.39
10	C	337	HEM	C3B-C4B	-8.59	1.44	1.51
10	C	340	HEM	C3B-C4B	-7.64	1.45	1.51
10	C	339	HEM	C2D-C3D	-6.56	1.34	1.54
10	C	338	HEM	C3B-C4B	-6.55	1.46	1.51
10	C	339	HEM	C3B-C4B	-6.41	1.46	1.51
7	L	304	BCB	C4D-CHA	-6.32	1.37	1.45
10	C	340	HEM	C2D-C3D	-6.28	1.35	1.54
10	C	337	HEM	C2D-C3D	-6.18	1.36	1.54
7	M	806	BCB	C4D-CHA	-6.07	1.37	1.45
10	C	338	HEM	C2D-C3D	-5.95	1.36	1.54
7	L	302	BCB	C4D-CHA	-5.89	1.37	1.45
10	C	339	HEM	C3D-C4D	-5.71	1.44	1.51
10	C	340	HEM	C3D-C4D	-5.67	1.44	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
10	C	338	HEM	C3D-C4D	-5.15	1.45	1.51
10	C	337	HEM	C3D-C4D	-4.81	1.45	1.51
9	M	501	MQ7	C11-C12	-4.63	1.43	1.50
10	C	337	HEM	C2C-C1C	-4.43	1.44	1.52
7	M	805	BCB	C4D-CHA	-4.38	1.39	1.45
10	C	340	HEM	C2C-C1C	-4.21	1.44	1.52
10	C	338	HEM	C2C-C1C	-4.14	1.44	1.52
12	L	707	LDA	CM1-N1	-4.09	1.43	1.49
7	L	302	BCB	O2D-CED	-4.07	1.35	1.45
10	C	339	HEM	C2C-C1C	-3.87	1.45	1.52
12	M	704	LDA	CM1-N1	-3.75	1.43	1.49
7	L	304	BCB	O2D-CED	-3.49	1.36	1.45
9	M	501	MQ7	C26-C27	-3.48	1.40	1.50
12	L	702	LDA	CM1-N1	-3.45	1.44	1.49
7	L	302	BCB	C1-C2	-3.38	1.37	1.49
8	L	402	BPB	O2D-CED	-3.27	1.37	1.45
12	H	701	LDA	CM1-N1	-3.05	1.44	1.49
6	M	802	SO4	O3-S	-3.02	1.36	1.47
12	M	706	LDA	C1-N1	-3.01	1.45	1.51
12	L	702	LDA	CM2-N1	-2.96	1.44	1.49
12	M	705	LDA	C1-N1	-2.96	1.46	1.51
12	H	701	LDA	CM2-N1	-2.89	1.45	1.49
12	M	705	LDA	CM1-N1	-2.84	1.45	1.49
7	M	806	BCB	CAA-CBA	-2.82	1.43	1.52
7	M	805	BCB	O2D-CED	-2.81	1.38	1.45
6	M	804	SO4	O1-S	-2.80	1.37	1.47
12	M	704	LDA	CM2-N1	-2.69	1.45	1.49
8	L	402	BPB	C4C-C3C	-2.67	1.39	1.45
12	M	704	LDA	C1-N1	-2.59	1.46	1.51
12	M	706	LDA	CM2-N1	-2.53	1.45	1.49
12	M	705	LDA	CM2-N1	-2.53	1.45	1.49
12	M	706	LDA	CM1-N1	-2.49	1.45	1.49
8	M	401	BPB	O2D-CED	-2.49	1.39	1.45
10	C	339	HEM	C2B-C1B	-2.46	1.43	1.51
12	H	703	LDA	CM1-N1	-2.43	1.45	1.49
8	M	401	BPB	C3B-C2B	-2.37	1.34	1.40
12	H	703	LDA	C1-N1	-2.36	1.47	1.51
9	M	501	MQ7	C10-C5	-2.34	1.37	1.40
10	C	340	HEM	C2D-C1D	-2.32	1.44	1.51
10	C	338	HEM	C2B-C1B	-2.20	1.44	1.51
10	C	337	HEM	C2B-C1B	-2.20	1.44	1.51
10	C	340	HEM	CAD-C3D	-2.16	1.49	1.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	M	802	SO4	O2-S	-2.14	1.39	1.47
9	M	501	MQ7	C41-C42	-2.10	1.44	1.50
10	C	340	HEM	C2B-C1B	-2.09	1.45	1.51
7	L	302	BCB	C4C-C3C	-2.08	1.41	1.45
10	C	338	HEM	C2D-C1D	-2.00	1.45	1.51
10	C	338	HEM	C1C-NC	2.00	1.38	1.36
7	M	806	BCB	C4C-NC	2.01	1.41	1.37
7	M	806	BCB	CHD-C4C	2.03	1.41	1.35
10	C	337	HEM	FE-NB	2.12	2.08	1.97
10	C	340	HEM	C1C-NC	2.13	1.38	1.36
10	C	337	HEM	C1C-NC	2.16	1.38	1.36
10	C	339	HEM	C4C-NC	2.18	1.38	1.36
7	M	805	BCB	C5-C3	2.31	1.56	1.51
10	C	338	HEM	C4C-NC	2.34	1.38	1.36
9	M	501	MQ7	C12-C13	2.40	1.37	1.33
8	L	402	BPB	C1A-CHA	2.41	1.41	1.36
9	M	501	MQ7	C42-C43	2.44	1.39	1.32
7	M	806	BCB	C1A-CHA	2.47	1.42	1.37
8	M	401	BPB	C1A-CHA	2.53	1.41	1.36
9	M	501	MQ7	C22-C23	2.57	1.38	1.33
8	L	402	BPB	CMD-C2D	2.62	1.56	1.50
8	M	401	BPB	O2D-CGD	2.64	1.39	1.33
7	L	302	BCB	C2-C3	2.87	1.38	1.33
10	C	339	HEM	C3B-CAB	2.89	1.56	1.51
9	M	501	MQ7	C27-C28	2.92	1.38	1.33
7	M	805	BCB	O2D-CGD	2.98	1.40	1.33
8	M	401	BPB	C3B-C4B	3.03	1.45	1.41
7	M	806	BCB	CAC-C3C	3.19	1.37	1.33
10	C	338	HEM	FE-NC	3.20	2.08	1.95
7	M	805	BCB	C2-C3	3.30	1.39	1.33
8	L	402	BPB	C2-C3	3.36	1.39	1.33
9	M	501	MQ7	C37-C38	3.40	1.39	1.33
7	M	806	BCB	C2-C3	3.43	1.39	1.33
9	M	501	MQ7	C17-C18	3.44	1.39	1.33
8	M	401	BPB	C2-C3	3.48	1.39	1.33
7	L	304	BCB	C2-C3	3.54	1.39	1.33
7	M	805	BCB	O2A-CGA	3.58	1.44	1.33
8	M	401	BPB	O2A-CGA	3.74	1.44	1.33
8	L	402	BPB	CAC-C3C	3.78	1.37	1.33
9	M	501	MQ7	C32-C33	3.78	1.40	1.33
10	C	338	HEM	CBC-CAC	3.91	1.51	1.29
10	C	340	HEM	CBB-CAB	3.96	1.52	1.29

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
10	C	338	HEM	CBB-CAB	3.99	1.52	1.29
8	L	402	BPB	O2D-CGD	4.01	1.43	1.33
10	C	337	HEM	CBB-CAB	4.11	1.53	1.29
8	L	402	BPB	C3B-C4B	4.17	1.46	1.41
10	C	337	HEM	CBC-CAC	4.27	1.54	1.29
10	C	340	HEM	CBC-CAC	4.38	1.54	1.29
7	M	805	BCB	CAC-C3C	4.49	1.38	1.33
10	C	339	HEM	CBB-CAB	4.52	1.55	1.29
10	C	339	HEM	CBC-CAC	4.56	1.55	1.29
7	L	302	BCB	CAC-C3C	4.90	1.39	1.33
8	M	401	BPB	CAC-C3C	5.15	1.39	1.33
7	L	304	BCB	CAC-C3C	5.31	1.39	1.33

All (116) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
10	C	338	HEM	C3C-CAC-CBC	-9.82	109.39	124.46
10	C	339	HEM	C3B-CAB-CBB	-8.88	110.84	124.46
10	C	337	HEM	C3B-CAB-CBB	-7.49	112.96	124.46
10	C	340	HEM	C3C-CAC-CBC	-7.20	113.41	124.46
10	C	338	HEM	C3B-CAB-CBB	-6.76	114.09	124.46
8	L	402	BPB	O1D-CGD-CBD	-6.74	114.96	124.62
8	M	401	BPB	O1D-CGD-CBD	-6.67	115.06	124.62
7	M	805	BCB	O1D-CGD-CBD	-6.66	115.08	124.62
7	L	304	BCB	O1D-CGD-CBD	-6.65	115.08	124.62
10	C	340	HEM	C3B-CAB-CBB	-6.22	114.92	124.46
10	C	337	HEM	C3C-CAC-CBC	-5.74	115.65	124.46
7	M	806	BCB	O1D-CGD-CBD	-5.63	116.56	124.62
10	C	339	HEM	C3C-CAC-CBC	-5.18	116.52	124.46
7	L	302	BCB	O1D-CGD-CBD	-5.05	117.38	124.62
7	M	806	BCB	OBD-CAD-CBD	-3.80	120.20	125.94
10	C	338	HEM	CBA-CAA-C2A	-3.66	105.98	112.53
7	M	806	BCB	CAA-C2A-C1A	-3.51	100.08	112.47
7	M	805	BCB	OBD-CAD-CBD	-3.49	120.67	125.94
7	L	304	BCB	OBD-CAD-CBD	-3.39	120.82	125.94
7	L	304	BCB	C4-C3-C5	-3.37	110.26	115.41
7	M	806	BCB	O2A-CGA-O1A	-3.33	114.90	123.49
9	M	501	MQ7	C34-C33-C35	-3.25	110.44	115.41
7	M	806	BCB	CBC-CAC-C3C	-3.19	119.70	127.07
8	M	401	BPB	C2C-C3C-C4C	-3.16	104.35	107.24
9	M	501	MQ7	C39-C38-C40	-3.09	110.68	115.41
7	L	302	BCB	OBD-CAD-CBD	-3.05	121.34	125.94

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
9	M	501	MQ7	C40-C41-C42	-3.00	103.84	111.69
7	M	805	BCB	C4-C3-C2	-2.67	118.27	123.50
8	L	402	BPB	CBD-CHA-C4D	-2.62	105.52	108.46
9	M	501	MQ7	C26-C25-C23	-2.57	104.33	112.71
12	H	701	LDA	CM2-N1-CM1	-2.54	105.97	108.83
11	M	600	NS5	C19-C20-C21	-2.54	123.53	127.20
8	L	402	BPB	CHD-C1D-ND	-2.50	119.98	124.66
7	L	302	BCB	O2A-CGA-O1A	-2.48	117.09	123.49
8	L	402	BPB	C2C-C3C-C4C	-2.41	105.04	107.24
7	L	302	BCB	C15-C13-C12	-2.40	97.94	112.27
10	C	337	HEM	CAA-C2A-C3A	-2.40	122.16	129.00
9	M	501	MQ7	C36-C35-C33	-2.33	105.13	112.71
7	M	805	BCB	OBB-CAB-C3B	-2.32	116.33	120.00
7	M	806	BCB	CBB-CAB-C3B	-2.32	113.46	120.33
10	C	340	HEM	CBA-CAA-C2A	-2.31	108.39	112.53
10	C	337	HEM	CBD-CAD-C3D	-2.26	106.99	113.55
11	M	600	NS5	C16-C15-C14	-2.25	114.36	118.10
8	M	401	BPB	CBD-CHA-C4D	-2.11	106.10	108.46
7	M	805	BCB	CBC-CAC-C3C	-2.09	122.24	127.07
11	M	600	NS5	C22-C21-C23	-2.02	114.73	118.10
7	L	304	BCB	C15-C13-C12	-2.01	100.27	112.27
7	M	806	BCB	C3C-C4C-NC	2.01	111.74	110.24
9	M	501	MQ7	C20-C21-C22	2.03	117.01	111.69
7	M	805	BCB	C4-C3-C5	2.04	118.52	115.41
8	L	402	BPB	CED-O2D-CGD	2.10	120.92	115.99
7	L	304	BCB	OBB-CAB-C3B	2.10	123.33	120.00
8	L	402	BPB	C2D-C1D-ND	2.19	113.01	109.73
8	M	401	BPB	OBB-CAB-C3B	2.19	123.47	120.00
9	M	501	MQ7	C35-C36-C37	2.22	117.51	111.69
12	H	701	LDA	O1-N1-CM1	2.29	112.12	109.05
7	L	304	BCB	O2A-CGA-CBA	2.32	118.95	111.90
8	M	401	BPB	CBD-CHA-C1A	2.33	130.71	126.78
6	M	802	SO4	O2-S-O1	2.38	117.05	109.50
8	L	402	BPB	O2A-CGA-CBA	2.42	119.26	111.90
7	M	805	BCB	C3C-C4C-NC	2.47	112.08	110.24
7	M	806	BCB	CMD-C2D-C3D	2.51	130.00	125.09
10	C	339	HEM	CMD-C2D-C3D	2.55	125.61	114.35
11	M	600	NS5	C18-C19-C20	2.60	129.13	123.39
7	L	304	BCB	CMB-C2B-C3B	2.60	130.18	125.09
8	M	401	BPB	C6-C5-C3	2.62	118.22	112.48
9	M	501	MQ7	C41-C42-C43	2.62	137.83	127.73
8	M	401	BPB	C3C-C2C-C1C	2.63	104.88	100.99

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	M	805	BCB	CMB-C2B-C3B	2.73	130.42	125.09
10	C	338	HEM	CMD-C2D-C3D	2.79	126.68	114.35
10	C	340	HEM	CMD-C2D-C3D	2.86	126.98	114.35
7	L	302	BCB	O2A-CGA-CBA	2.99	121.00	111.90
10	C	337	HEM	CAA-C2A-C1A	3.06	130.33	127.01
8	L	402	BPB	C3C-C2C-C1C	3.06	105.51	100.99
7	L	304	BCB	C3C-C4C-NC	3.09	112.54	110.24
10	C	337	HEM	C2D-C3D-C4D	3.10	106.75	101.50
9	M	501	MQ7	C25-C23-C22	3.10	126.93	121.05
8	M	401	BPB	CED-O2D-CGD	3.12	123.32	115.99
10	C	340	HEM	CAD-C3D-C4D	3.14	123.53	112.47
7	L	302	BCB	CMB-C2B-C3B	3.19	131.34	125.09
10	C	339	HEM	CAD-C3D-C4D	3.22	123.82	112.47
7	M	805	BCB	C6-C5-C3	3.27	119.66	112.48
10	C	338	HEM	C2D-C3D-C4D	3.27	107.04	101.50
10	C	339	HEM	C2D-C3D-C4D	3.28	107.06	101.50
7	L	302	BCB	C3C-C4C-NC	3.29	112.69	110.24
11	M	600	NS5	C19-C18-C17	3.33	130.76	123.39
7	M	805	BCB	C4A-NA-C1A	3.41	110.06	106.04
7	L	304	BCB	C4A-NA-C1A	3.41	110.06	106.04
9	M	501	MQ7	C35-C33-C32	3.45	127.59	121.05
10	C	340	HEM	C2D-C3D-C4D	3.49	107.41	101.50
10	C	338	HEM	CAD-C3D-C4D	3.53	124.92	112.47
10	C	337	HEM	CAD-C3D-C2D	3.56	123.45	113.22
10	C	337	HEM	CMB-C2B-C3B	3.67	125.70	116.53
7	L	302	BCB	C4A-NA-C1A	3.81	110.53	106.04
10	C	339	HEM	CMC-C2C-C3C	3.84	126.11	116.53
7	M	806	BCB	C4A-NA-C1A	4.16	110.94	106.04
7	M	806	BCB	O2A-CGA-CBA	4.26	124.89	111.90
7	L	302	BCB	OBB-CAB-C3B	4.34	126.88	120.00
10	C	340	HEM	CMC-C2C-C3C	4.39	127.50	116.53
10	C	338	HEM	CMB-C2B-C3B	4.48	127.71	116.53
10	C	338	HEM	CMC-C2C-C3C	4.53	127.84	116.53
10	C	337	HEM	CAD-C3D-C4D	4.91	129.80	112.47
10	C	338	HEM	CAD-C3D-C2D	5.10	127.88	113.22
7	M	806	BCB	CMB-C2B-C3B	5.23	135.32	125.09
10	C	340	HEM	CAD-C3D-C2D	5.30	128.47	113.22
10	C	339	HEM	CAD-C3D-C2D	5.53	129.11	113.22
10	C	340	HEM	CMB-C2B-C3B	5.53	130.34	116.53
10	C	337	HEM	CMC-C2C-C3C	5.57	130.44	116.53
7	M	806	BCB	OBB-CAB-C3B	5.76	129.12	120.00
7	L	302	BCB	O2D-CGD-CBD	5.88	119.36	111.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
10	C	339	HEM	CMB-C2B-C3B	6.19	131.99	116.53
7	M	806	BCB	O2D-CGD-CBD	7.90	122.14	111.30
8	L	402	BPB	O2D-CGD-CBD	8.22	122.57	111.30
7	M	805	BCB	O2D-CGD-CBD	8.72	123.26	111.30
7	L	304	BCB	O2D-CGD-CBD	9.00	123.64	111.30
8	M	401	BPB	O2D-CGD-CBD	9.38	124.16	111.30

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

10 monomers are involved in 49 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
10	C	339	HEM	1	0
7	L	302	BCB	5	0
7	L	304	BCB	4	0
8	L	402	BPB	5	0
12	L	702	LDA	8	0
12	L	707	LDA	2	0
8	M	401	BPB	12	0
12	M	706	LDA	2	0
7	M	805	BCB	7	0
7	M	806	BCB	6	0

## 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, 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	C	332/336 (98%)	-0.84	1 (0%) 94 94	11, 25, 47, 67	15 (4%)
2	L	273/273 (100%)	-0.99	1 (0%) 93 93	9, 19, 41, 50	7 (2%)
3	M	323/323 (100%)	-0.95	1 (0%) 94 94	9, 22, 50, 61	8 (2%)
4	H	249/258 (96%)	-0.80	0 100 100	12, 30, 54, 64	22 (8%)
All	All	1177/1190 (98%)	-0.89	3 (0%) 94 94	9, 24, 49, 67	52 (4%)

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	L	59	TRP	2.3
3	M	319	PRO	2.2
1	C	332	LYS	2.0

### 6.2 Non-standard residues in protein, DNA, RNA chains [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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
4	FME	H	1[B]	10/11	0.97	0.09	-	27,30,30,30	10
4	FME	H	1[A]	10/11	0.97	0.09	-	22,25,27,27	10

### 6.3 Carbohydrates [i](#)

There are no carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
12	LDA	L	707	16/16	0.64	0.27	9.77	54,64,65,66	0
12	LDA	L	702	16/16	0.80	0.25	6.48	19,44,66,66	0
11	NS5	M	600	40/40	0.88	0.19	4.31	31,49,68,68	9
12	LDA	M	704	16/16	0.83	0.22	3.54	53,59,61,62	3
12	LDA	M	705	16/16	0.90	0.15	3.12	51,55,59,59	5
12	LDA	M	706	16/16	0.74	0.27	2.06	69,70,72,72	4
9	MQ7	M	501	48/48	0.95	0.11	1.73	11,17,43,49	0
12	LDA	H	703	16/16	0.94	0.16	1.64	30,41,52,52	1
12	LDA	H	701	16/16	0.96	0.10	1.37	28,30,31,32	0
7	BCB	M	805	66/66	0.96	0.12	0.94	8,18,37,38	20
6	SO4	H	801	5/5	0.98	0.09	0.93	64,65,65,67	0
6	SO4	M	802	5/5	0.99	0.08	0.86	30,30,32,39	0
6	SO4	M	804	5/5	0.99	0.14	0.78	40,41,42,43	0
8	BPB	M	401	65/65	0.97	0.10	0.69	10,24,61,62	7
7	BCB	L	302	66/66	0.99	0.10	0.56	7,11,18,19	0
7	BCB	M	806	66/66	0.98	0.10	0.46	4,12,30,31	0
10	HEM	C	338	43/43	0.98	0.13	0.46	16,25,32,38	0
7	BCB	L	304	66/66	0.98	0.12	0.36	6,13,26,36	0
8	BPB	L	402	65/65	0.98	0.08	0.26	7,12,18,20	0
10	HEM	C	340	43/43	0.98	0.10	0.26	11,19,34,42	0
10	HEM	C	337	43/43	0.98	0.08	-0.59	21,27,38,43	0
10	HEM	C	339	43/43	0.99	0.06	-0.74	8,14,18,25	0
5	FE2	M	500	1/1	1.00	0.04	-2.10	14,14,14,14	0
6	SO4	M	803	5/5	0.96	0.11	-	85,85,86,86	0

## 6.5 Other polymers

There are no such residues in this entry.