



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 05:12 PM UTC

PDB ID : 5FOI / pdb_00005foi
Title : Crystal structure of mycinamicin VIII C21 methyl hydroxylase MycCI from *Micromonospora griseorubida* bound to mycinamicin VIII
Authors : Demars, M.; Sheng, F.; Podust, L.M.; Sherman, D.H.
Deposited on : 2015-11-22
Resolution : 2.21 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

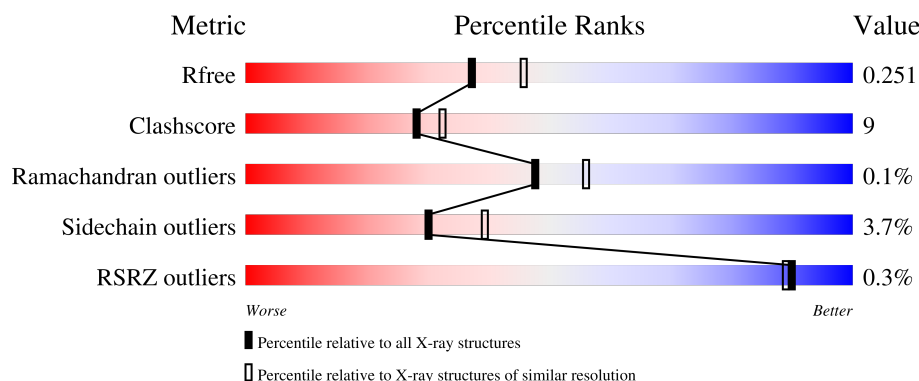
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

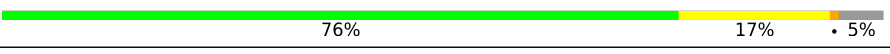
The reported resolution of this entry is 2.21 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	7682 (2.24-2.20)
Clashscore	190562	8402 (2.24-2.20)
Ramachandran outliers	187476	8303 (2.24-2.20)
Sidechain outliers	187428	8304 (2.24-2.20)
RSRZ outliers	180081	7683 (2.24-2.20)

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

Mol	Chain	Length	Quality of chain
1	A	408	 75% 19% • 5%
1	B	408	 76% 17% • 5%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 6369 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 MYCINAMICIN VIII C21 METHYL HYDROXYLASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	389	Total	C	N	O	S	0	0	0
			3004	1903	542	547	12			
1	B	388	Total	C	N	O	S	0	1	0
			3018	1909	549	547	13			

There are 40 discrepancies between the modelled and reference sequences:

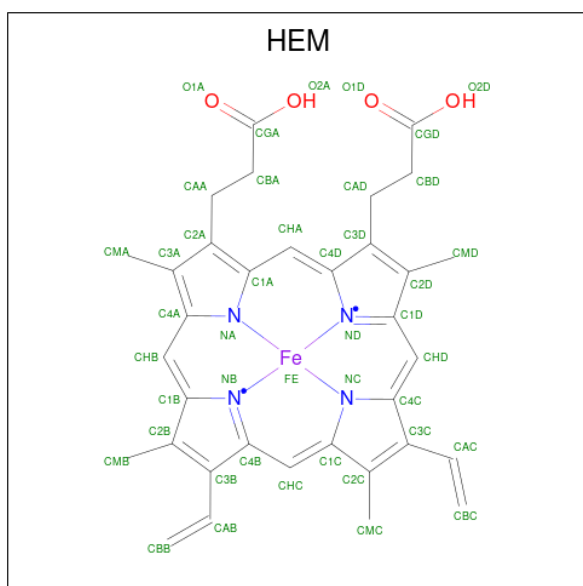
Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	expression tag	UNP Q83WF5
A	-18	GLY	-	expression tag	UNP Q83WF5
A	-17	SER	-	expression tag	UNP Q83WF5
A	-16	SER	-	expression tag	UNP Q83WF5
A	-15	HIS	-	expression tag	UNP Q83WF5
A	-14	HIS	-	expression tag	UNP Q83WF5
A	-13	HIS	-	expression tag	UNP Q83WF5
A	-12	HIS	-	expression tag	UNP Q83WF5
A	-11	HIS	-	expression tag	UNP Q83WF5
A	-10	HIS	-	expression tag	UNP Q83WF5
A	-9	SER	-	expression tag	UNP Q83WF5
A	-8	SER	-	expression tag	UNP Q83WF5
A	-7	GLY	-	expression tag	UNP Q83WF5
A	-6	LEU	-	expression tag	UNP Q83WF5
A	-5	VAL	-	expression tag	UNP Q83WF5
A	-4	PRO	-	expression tag	UNP Q83WF5
A	-3	ARG	-	expression tag	UNP Q83WF5
A	-2	GLY	-	expression tag	UNP Q83WF5
A	-1	SER	-	expression tag	UNP Q83WF5
A	0	HIS	-	expression tag	UNP Q83WF5
B	-19	MET	-	expression tag	UNP Q83WF5
B	-18	GLY	-	expression tag	UNP Q83WF5
B	-17	SER	-	expression tag	UNP Q83WF5
B	-16	SER	-	expression tag	UNP Q83WF5
B	-15	HIS	-	expression tag	UNP Q83WF5

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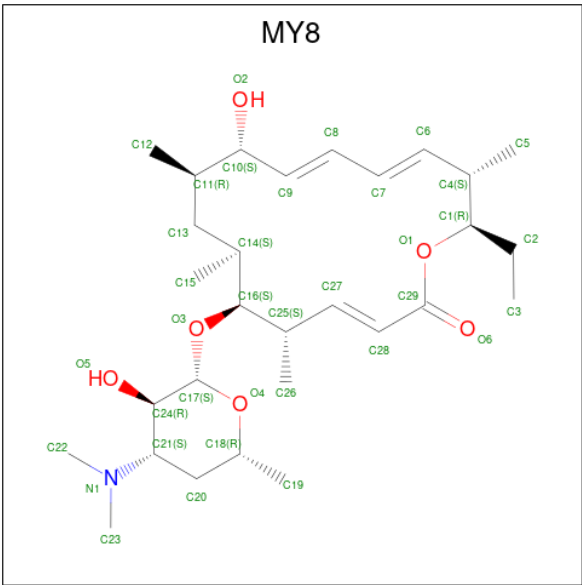
Chain	Residue	Modelled	Actual	Comment	Reference
B	-14	HIS	-	expression tag	UNP Q83WF5
B	-13	HIS	-	expression tag	UNP Q83WF5
B	-12	HIS	-	expression tag	UNP Q83WF5
B	-11	HIS	-	expression tag	UNP Q83WF5
B	-10	HIS	-	expression tag	UNP Q83WF5
B	-9	SER	-	expression tag	UNP Q83WF5
B	-8	SER	-	expression tag	UNP Q83WF5
B	-7	GLY	-	expression tag	UNP Q83WF5
B	-6	LEU	-	expression tag	UNP Q83WF5
B	-5	VAL	-	expression tag	UNP Q83WF5
B	-4	PRO	-	expression tag	UNP Q83WF5
B	-3	ARG	-	expression tag	UNP Q83WF5
B	-2	GLY	-	expression tag	UNP Q83WF5
B	-1	SER	-	expression tag	UNP Q83WF5
B	0	HIS	-	expression tag	UNP Q83WF5

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



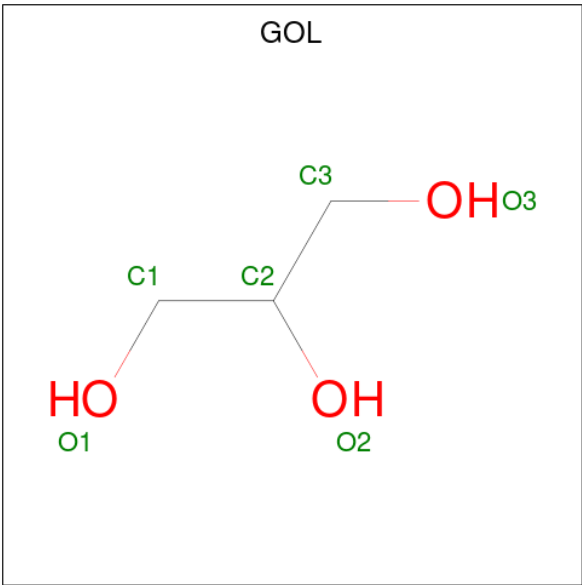
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
2	B	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		

- Molecule 3 is Mycinamicin VIII (CCD ID: MY8) (formula: $C_{29}H_{49}NO_6$).



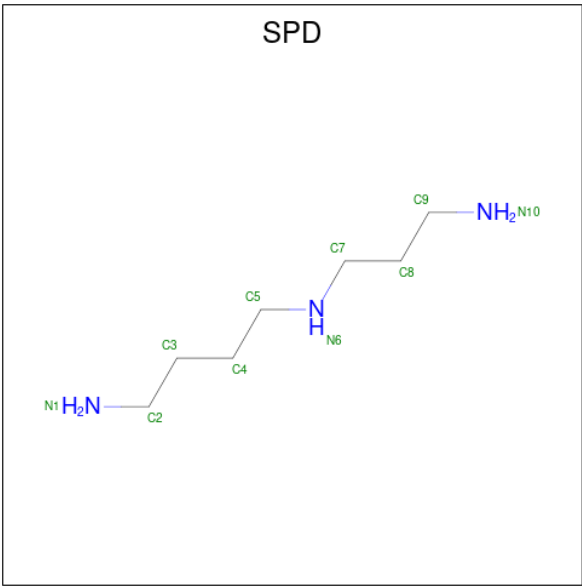
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			36	29	1	6		
3	B	1	Total	C	N	O	0	0
			36	29	1	6		

- Molecule 4 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



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

- Molecule 5 is SPERMIDINE (CCD ID: SPD) (formula: C₇H₁₉N₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	N	0	0
			10	7	3		
5	A	1	Total	C	N	0	0
			10	7	3		
5	A	1	Total	C	N	0	0
			10	7	3		

- Molecule 6 is water.

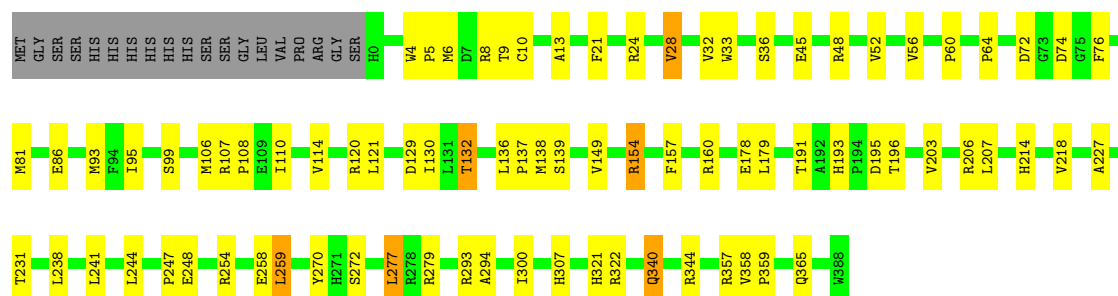
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	77	Total	O	0	0
			77	77		
6	B	70	Total	O	0	0
			70	70		

3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry 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.

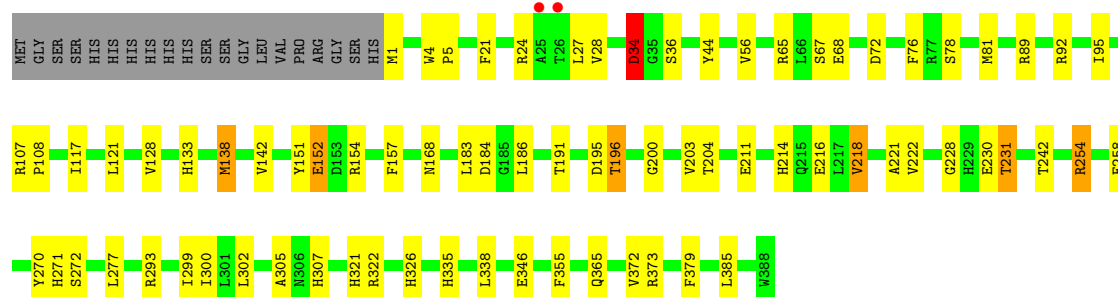
• Molecule 1: MYCINAMICIN VIII C21 METHYL HYDROXYLASE

Chain A: 



• Molecule 1: MYCINAMICIN VIII C21 METHYL HYDROXYLASE

Chain B: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	54.78Å 59.52Å 74.62Å 83.20° 72.15° 62.61°	Depositor
Resolution (Å)	71.00 – 2.21 71.00 – 2.21	Depositor EDS
% Data completeness (in resolution range)	90.8 (71.00-2.21) 90.8 (71.00-2.21)	Depositor EDS
R_{merge}	0.24	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.60 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.182 , 0.252 0.189 , 0.251	Depositor DCC
R_{free} test set	1838 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	26.0	Xtriage
Anisotropy	0.500	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 33.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.010 for -h,-h+k,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6369	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.79% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: SPD, HEM, MY8, GOL

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	1.02	0/3070	1.09	11/4188 (0.3%)
1	B	0.97	0/3083	1.07	7/4201 (0.2%)
All	All	0.99	0/6153	1.08	18/8389 (0.2%)

There are no bond length outliers.

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	24	ARG	N-CA-C	8.53	120.90	110.41
1	A	193	HIS	CA-C-N	-6.56	113.01	119.76
1	A	193	HIS	C-N-CA	-6.56	113.01	119.76
1	B	196	THR	N-CA-C	-5.91	101.21	110.36
1	A	28	VAL	CB-CA-C	-5.87	105.48	110.94
1	A	95	ILE	N-CA-C	5.73	116.47	110.62
1	B	322	ARG	CA-C-N	-5.70	113.37	119.98
1	B	322	ARG	C-N-CA	-5.70	113.37	119.98
1	A	130	ILE	N-CA-C	5.51	116.24	110.62
1	B	117	ILE	N-CA-C	-5.40	105.45	110.53
1	A	340	GLN	N-CA-C	5.39	117.91	111.71
1	A	139	SER	N-CA-C	-5.38	105.50	111.36
1	B	228	GLY	N-CA-C	5.22	119.00	112.73
1	A	178	GLU	N-CA-C	5.18	116.93	111.28
1	A	149	VAL	CA-C-N	-5.18	113.86	120.23
1	A	149	VAL	C-N-CA	-5.18	113.86	120.23
1	B	34	ASP	N-CA-C	5.11	118.61	112.38
1	B	218	VAL	CB-CA-C	-5.08	105.47	111.97

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3004	0	2974	45	0
1	B	3018	0	3004	58	0
2	A	43	0	30	2	0
2	B	43	0	30	8	0
3	A	36	0	0	1	0
3	B	36	0	0	0	0
4	A	12	0	16	3	0
5	A	30	0	57	3	0
6	A	77	0	0	5	0
6	B	70	0	0	2	0
All	All	6369	0	6111	110	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 9.

All (110) 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:B:78:SER:H	1:B:81:MET:HE3	1.50	0.76
1:B:76:PHE:CE2	1:B:81:MET:HE1	2.22	0.74
1:A:129:ASP:OD2	1:A:132:THR:OG1	2.06	0.73
1:A:340:GLN:HE21	1:A:344:ARG:HH12	1.38	0.70
1:B:138:MET:CE	1:B:142:VAL:HG21	2.24	0.68
1:B:195:ASP:OD2	1:B:196:THR:O	2.11	0.68
1:B:231:THR:HG21	2:B:413:HEM:CHC	2.25	0.65
1:B:365:GLN:HE21	1:B:385:LEU:H	1.46	0.64
1:A:4:TRP:CG	1:A:5:PRO:HA	2.33	0.64
1:B:76:PHE:HE2	1:B:81:MET:HE1	1.61	0.63
1:B:365:GLN:NE2	1:B:385:LEU:H	1.97	0.62
1:B:231:THR:HG21	2:B:413:HEM:C4B	2.35	0.61
1:B:121:LEU:HD11	1:B:128:VAL:HG11	1.80	0.61
1:B:4:TRP:CD1	1:B:299:ILE:HD13	2.37	0.59
1:A:4:TRP:CD2	1:A:5:PRO:HA	2.38	0.59
1:B:21:PHE:CG	1:B:28:VAL:HG21	2.38	0.59
1:B:107:ARG:HB3	1:B:108:PRO:HD3	1.85	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:200:GLY:O	1:B:204:THR:HG22	2.02	0.58
1:B:34:ASP:HB3	1:B:36:SER:H	1.69	0.57
1:B:89:ARG:NH1	1:B:211:GLU:O	2.38	0.57
1:B:65:ARG:NH2	1:B:68:GLU:O	2.38	0.57
1:A:136:LEU:HB3	1:A:137:PRO:HD3	1.87	0.56
1:B:1:MET:HA	6:B:2002:HOH:O	2.04	0.56
6:A:2054:HOH:O	1:B:326:HIS:CD2	2.57	0.56
1:A:107:ARG:HB3	1:A:108:PRO:HD3	1.86	0.56
5:A:1391:SPD:H42	5:A:1392:SPD:H92	1.87	0.56
2:A:413:HEM:HMB2	2:A:413:HEM:HBB2	1.88	0.55
1:A:36:SER:HB2	1:A:64:PRO:HB3	1.89	0.54
1:A:227:ALA:O	1:A:231:THR:OG1	2.25	0.54
1:A:254:ARG:HH11	1:A:359:PRO:CB	2.20	0.53
1:B:254:ARG:HG2	1:B:254:ARG:HH11	1.75	0.52
1:A:277:LEU:HB2	1:A:300:ILE:HB	1.92	0.52
1:A:106:MET:O	1:A:110:ILE:HG13	2.10	0.51
2:B:413:HEM:HBB2	2:B:413:HEM:HMB2	1.91	0.51
2:A:413:HEM:HBB2	2:A:413:HEM:CMB	2.40	0.51
1:B:138:MET:CE	1:B:142:VAL:CG2	2.89	0.51
1:B:271:HIS:O	1:B:272:SER:C	2.53	0.51
1:B:231:THR:CG2	2:B:413:HEM:CHC	2.89	0.50
5:A:1391:SPD:C4	5:A:1392:SPD:H92	2.41	0.50
1:A:207:LEU:C	1:A:207:LEU:HD23	2.37	0.50
1:A:32:VAL:O	1:A:33:TRP:C	2.51	0.50
1:A:247:PRO:O	1:A:248:GLU:C	2.54	0.50
1:A:258:GLU:CB	4:A:1390:GOL:C2	2.90	0.50
1:B:231:THR:HG23	2:B:413:HEM:CAB	2.42	0.49
1:A:259:LEU:HD21	1:A:321:HIS:CE1	2.47	0.49
1:A:10:CYS:HB3	1:A:13:ALA:HB3	1.95	0.48
1:A:56:VAL:HG13	1:A:279:ARG:HB3	1.94	0.48
1:A:258:GLU:CB	4:A:1390:GOL:H2	2.43	0.48
1:B:4:TRP:CG	1:B:299:ILE:HD13	2.48	0.47
1:B:254:ARG:HD3	1:B:355:PHE:O	2.13	0.47
1:B:138:MET:HE1	1:B:346:GLU:CD	2.39	0.47
1:B:138:MET:O	1:B:138:MET:HE2	2.14	0.47
1:A:160:ARG:HB3	1:A:179:LEU:HG	1.97	0.47
1:B:89:ARG:NH2	1:B:216:GLU:OE1	2.43	0.47
1:B:184:ASP:OD2	1:B:214:HIS:HE1	1.98	0.47
1:A:293:ARG:O	1:A:294:ALA:C	2.58	0.47
1:B:95:ILE:HA	1:B:338:LEU:HD11	1.97	0.47
1:B:372:VAL:HG13	1:B:379:PHE:HE1	1.80	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:6:MET:HE3	1:A:6:MET:HB2	1.81	0.46
1:B:270:TYR:O	1:B:307:HIS:HE1	1.97	0.46
1:A:21:PHE:CD1	1:A:28:VAL:HG21	2.50	0.46
1:B:191:THR:HA	1:B:203:VAL:HG11	1.98	0.45
1:B:67:SER:OG	1:B:168:ASN:ND2	2.49	0.45
1:A:322:ARG:O	1:A:322:ARG:CG	2.65	0.45
1:A:357:ARG:O	1:A:359:PRO:HD3	2.16	0.45
1:B:121:LEU:HD13	1:B:133:HIS:HB3	1.99	0.45
3:A:414:MY8:C17	3:A:414:MY8:C26	2.95	0.44
1:A:191:THR:HA	1:A:203:VAL:HG11	1.99	0.44
1:B:21:PHE:CD1	1:B:28:VAL:HG21	2.51	0.44
1:B:231:THR:CG2	2:B:413:HEM:C4B	3.00	0.44
1:A:48:ARG:O	1:A:52:VAL:HG22	2.17	0.44
1:B:230:GLU:OE2	1:B:373:ARG:NH1	2.49	0.44
2:B:413:HEM:HBB2	2:B:413:HEM:CMB	2.48	0.44
1:B:89:ARG:NH2	1:B:216:GLU:OE2	2.51	0.43
5:A:1391:SPD:C4	5:A:1392:SPD:C9	2.97	0.43
1:B:218:VAL:O	1:B:222:VAL:HG23	2.18	0.43
1:A:45:GLU:HA	6:A:2015:HOH:O	2.18	0.43
1:B:4:TRP:CD2	1:B:5:PRO:HA	2.54	0.43
1:A:254:ARG:NH1	1:A:359:PRO:CB	2.82	0.43
1:B:335:HIS:HD2	2:B:413:HEM:O1D	2.02	0.43
1:A:93:MET:HG2	1:A:206:ARG:HD2	2.00	0.43
1:A:244:LEU:HD23	1:A:244:LEU:HA	1.82	0.43
1:A:8:ARG:NH2	1:A:272:SER:O	2.49	0.43
1:B:78:SER:N	1:B:81:MET:HE3	2.28	0.42
1:B:151:TYR:CE1	1:B:154:ARG:HD3	2.54	0.42
1:B:277:LEU:HD11	1:B:302:LEU:HD11	2.01	0.42
1:A:60:PRO:CG	1:A:81:MET:HE2	2.50	0.42
1:B:44:TYR:HA	1:B:305:ALA:HB1	2.01	0.42
1:B:258:GLU:HG2	6:B:2055:HOH:O	2.20	0.42
1:B:326:HIS:CD2	1:B:326:HIS:H	2.37	0.42
1:A:270:TYR:O	1:A:307:HIS:HE1	2.02	0.42
1:B:277:LEU:HB2	1:B:300:ILE:HB	2.00	0.42
1:A:120:ARG:NH2	6:A:2040:HOH:O	2.53	0.42
1:A:195:ASP:OD2	1:A:196:THR:O	2.37	0.42
1:B:152:GLU:H	1:B:152:GLU:CD	2.28	0.42
1:B:138:MET:HE2	1:B:142:VAL:CG2	2.49	0.42
1:A:258:GLU:CB	4:A:1390:GOL:O2	2.68	0.41
1:A:120:ARG:NH1	6:A:2042:HOH:O	2.39	0.41
1:A:74:ASP:HA	6:A:2026:HOH:O	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:214:HIS:O	1:A:218:VAL:HG23	2.21	0.41
1:A:99:SER:OG	1:B:321:HIS:HB3	2.21	0.41
1:A:238:LEU:O	1:A:241:LEU:HB3	2.21	0.41
1:A:322:ARG:O	1:A:322:ARG:HG3	2.21	0.41
1:B:242:THR:OG1	1:B:271:HIS:CE1	2.74	0.41
1:A:76:PHE:CD1	1:A:76:PHE:C	2.99	0.41
1:B:154:ARG:HA	1:B:157:PHE:CE2	2.56	0.41
1:B:254:ARG:HH11	1:B:254:ARG:CG	2.34	0.40
1:A:154:ARG:HA	1:A:157:PHE:CE2	2.55	0.40
1:B:183:LEU:HD22	1:B:221:ALA:CB	2.52	0.40
1:B:183:LEU:HD22	1:B:221:ALA:HB1	2.03	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	387/408 (95%)	369 (95%)	18 (5%)	0	100	100
1	B	387/408 (95%)	372 (96%)	14 (4%)	1 (0%)	36	41
All	All	774/816 (95%)	741 (96%)	32 (4%)	1 (0%)	48	56

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	27	LEU

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	313/347 (90%)	301 (96%)	12 (4%)	29	38
1	B	316/347 (91%)	305 (96%)	11 (4%)	32	41
All	All	629/694 (91%)	606 (96%)	23 (4%)	30	39

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

Mol	Chain	Res	Type
1	A	9	THR
1	A	72	ASP
1	A	86	GLU
1	A	114	VAL
1	A	121	LEU
1	A	132	THR
1	A	138	MET
1	A	154	ARG
1	A	259	LEU
1	A	277	LEU
1	A	358	VAL
1	A	365	GLN
1	B	24	ARG
1	B	34	ASP
1	B	56	VAL
1	B	72	ASP
1	B	92	ARG
1	B	138	MET
1	B	152	GLU
1	B	186	LEU
1	B	231	THR
1	B	254	ARG
1	B	293	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	133	HIS
1	A	141	GLN
1	A	307	HIS
1	A	321	HIS

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Mol	Chain	Res	Type
1	A	336	GLN
1	A	340	GLN
1	B	141	GLN
1	B	168	ASN
1	B	214	HIS
1	B	215	GLN
1	B	271	HIS
1	B	307	HIS
1	B	326	HIS
1	B	365	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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

9 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	$\# Z > 2$
4	GOL	A	1389	-	5,5,5	0.29	0	5,5,5	0.68	0
5	SPD	A	1392	-	9,9,9	0.40	0	8,8,8	1.22	0
5	SPD	A	1393	-	9,9,9	0.27	0	8,8,8	0.98	0
3	MY8	A	414	-	36,37,37	2.18	4 (11%)	43,51,51	4.03	25 (58%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	SPD	A	1391	-	9,9,9	0.42	0	8,8,8	1.28	2 (25%)
4	GOL	A	1390	-	5,5,5	0.57	0	5,5,5	0.70	0
2	HEM	B	413	1	50,50,50	1.69	13 (26%)	67,82,82	2.07	21 (31%)
3	MY8	B	414	-	36,37,37	2.24	4 (11%)	43,51,51	3.62	23 (53%)
2	HEM	A	413	1	50,50,50	1.67	9 (18%)	67,82,82	1.93	17 (25%)

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	GOL	A	1389	-	-	0/4/4/4	-
5	SPD	A	1392	-	-	5/7/7/7	-
5	SPD	A	1393	-	-	1/7/7/7	-
3	MY8	A	414	-	-	9/46/62/62	0/1/2/2
5	SPD	A	1391	-	-	4/7/7/7	-
4	GOL	A	1390	-	-	2/4/4/4	-
2	HEM	B	413	1	-	1/14/54/54	-
3	MY8	B	414	-	-	9/46/62/62	0/1/2/2
2	HEM	A	413	1	-	0/14/54/54	-

All (30) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	414	MY8	O2-C10	-11.59	1.22	1.43
3	A	414	MY8	O2-C10	-11.50	1.22	1.43
2	A	413	HEM	FE-NC	5.06	2.11	1.95
2	A	413	HEM	FE-NB	4.83	2.09	1.94
2	B	413	HEM	FE-NC	4.31	2.09	1.95
2	B	413	HEM	FE-NB	3.92	2.07	1.94
2	A	413	HEM	C4B-NB	-3.86	1.31	1.38
3	B	414	MY8	O1-C1	-3.34	1.40	1.46
2	A	413	HEM	C1B-NB	-3.25	1.34	1.40
2	B	413	HEM	C4B-NB	-3.23	1.32	1.38
3	A	414	MY8	C14-C16	-3.02	1.48	1.53
2	B	413	HEM	C1B-NB	-2.90	1.35	1.40
2	B	413	HEM	C4D-ND	-2.80	1.35	1.40
2	B	413	HEM	C1D-ND	-2.71	1.33	1.38
2	B	413	HEM	C1C-C2C	-2.65	1.40	1.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	414	MY8	C13-C11	-2.60	1.50	1.54
2	B	413	HEM	FE-NA	2.57	2.03	1.95
2	A	413	HEM	C1C-NC	-2.46	1.35	1.39
2	A	413	HEM	C1D-C2D	2.40	1.49	1.44
3	A	414	MY8	C25-C16	-2.34	1.48	1.54
2	B	413	HEM	C1A-C2A	-2.31	1.40	1.44
2	B	413	HEM	FE-ND	-2.31	1.87	1.94
2	A	413	HEM	C4D-C3D	2.27	1.48	1.45
2	B	413	HEM	O2D-CGD	-2.24	1.23	1.30
2	B	413	HEM	CHB-C1B	2.21	1.42	1.38
3	B	414	MY8	C14-C16	-2.19	1.50	1.53
2	A	413	HEM	O1A-CGA	2.18	1.29	1.22
3	B	414	MY8	O1-C29	2.10	1.39	1.34
2	B	413	HEM	C4C-NC	-2.10	1.35	1.39
2	A	413	HEM	C1C-C2C	-2.05	1.41	1.45

All (88) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	414	MY8	C15-C14-C16	-9.61	95.46	111.14
3	A	414	MY8	C5-C4-C1	-9.43	97.98	111.78
3	B	414	MY8	C5-C4-C1	-8.32	99.61	111.78
3	A	414	MY8	C5-C4-C6	-8.14	91.49	109.91
3	A	414	MY8	C10-C9-C8	-7.80	111.46	124.88
3	A	414	MY8	C1-O1-C29	-7.71	106.13	117.50
3	A	414	MY8	C12-C11-C13	-6.90	100.23	110.68
3	B	414	MY8	O1-C1-C2	-6.63	96.33	106.90
3	A	414	MY8	C26-C25-C16	-6.33	99.98	111.11
3	A	414	MY8	O4-C18-C20	6.21	118.22	109.34
3	B	414	MY8	C26-C25-C27	-6.13	96.03	109.91
3	B	414	MY8	O4-C17-C24	-6.02	98.00	110.37
3	A	414	MY8	O1-C1-C2	-5.93	97.46	106.90
2	A	413	HEM	CHC-C4B-NB	5.89	130.77	124.42
3	B	414	MY8	C1-O1-C29	-5.78	108.98	117.50
2	B	413	HEM	CHC-C4B-NB	5.68	130.54	124.42
3	A	414	MY8	O4-C17-C24	-5.65	98.76	110.37
2	B	413	HEM	CHA-C4D-ND	5.61	131.30	124.37
3	A	414	MY8	C18-C20-C21	-5.54	101.01	110.46
3	B	414	MY8	C10-C9-C8	-5.43	115.54	124.88
3	B	414	MY8	C5-C4-C6	-5.17	98.20	109.91
2	B	413	HEM	C1B-NB-C4B	5.17	111.33	105.21
2	A	413	HEM	C1B-NB-C4B	4.84	110.94	105.21

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	413	HEM	CHD-C1D-ND	4.70	129.48	124.42
3	A	414	MY8	C26-C25-C27	-4.63	99.44	109.91
3	B	414	MY8	C18-C20-C21	-4.60	102.61	110.46
3	B	414	MY8	O4-C18-C20	4.52	115.81	109.34
3	A	414	MY8	C15-C14-C13	-4.44	103.95	110.68
2	B	413	HEM	CHA-C4D-C3D	-4.43	117.06	125.23
2	B	413	HEM	CHD-C1D-ND	4.19	128.94	124.42
2	B	413	HEM	CAC-C3C-C4C	4.09	134.60	124.82
3	B	414	MY8	O4-C18-C19	4.03	114.43	106.88
3	A	414	MY8	C15-C14-C16	-4.03	104.56	111.14
3	B	414	MY8	C17-C24-C21	3.97	115.52	109.15
3	B	414	MY8	C20-C21-C24	-3.92	104.38	110.02
3	A	414	MY8	C4-C6-C7	-3.89	118.06	126.11
2	A	413	HEM	CMC-C2C-C1C	3.81	131.43	124.73
3	A	414	MY8	O3-C16-C25	-3.77	103.35	110.93
3	A	414	MY8	C17-C24-C21	3.77	115.19	109.15
3	B	414	MY8	O2-C10-C9	3.75	120.57	110.88
2	A	413	HEM	CAC-C3C-C4C	3.55	133.30	124.82
3	A	414	MY8	O2-C10-C9	3.51	119.94	110.88
3	B	414	MY8	C4-C6-C7	-3.48	118.91	126.11
3	A	414	MY8	O3-C17-O4	-3.47	101.57	110.69
3	B	414	MY8	C13-C14-C16	3.33	119.72	111.58
2	B	413	HEM	C3B-C4B-NB	-3.31	107.09	109.47
2	A	413	HEM	CHA-C4D-ND	3.26	128.40	124.37
2	B	413	HEM	CAC-C3C-C2C	-3.24	117.90	128.43
2	A	413	HEM	CAC-C3C-C2C	-3.11	118.31	128.43
2	A	413	HEM	C3B-C4B-NB	-3.10	107.24	109.47
3	B	414	MY8	O1-C29-O6	3.09	128.28	123.34
3	A	414	MY8	O3-C17-C24	3.06	115.63	108.09
3	A	414	MY8	C17-O3-C16	-3.05	110.76	117.98
2	A	413	HEM	CAD-C3D-C4D	2.97	129.87	124.70
2	B	413	HEM	O2D-CGD-O1D	-2.93	115.81	123.33
2	A	413	HEM	CHD-C1D-C2D	-2.73	120.72	125.03
2	B	413	HEM	C3B-C2B-C1B	-2.72	104.37	106.41
2	B	413	HEM	CHD-C1D-C2D	-2.70	120.76	125.03
2	B	413	HEM	CMB-C2B-C1B	2.69	129.24	125.03
2	B	413	HEM	C4A-C3A-C2A	2.69	109.89	106.82
2	B	413	HEM	O2A-CGA-CBA	2.65	122.37	114.00
2	B	413	HEM	CHD-C4C-C3C	2.63	129.64	125.21
3	A	414	MY8	C19-C18-C20	2.62	117.23	113.27
2	B	413	HEM	O2D-CGD-CBD	2.59	122.17	114.00
5	A	1391	SPD	C4-C5-N6	-2.57	105.16	112.07

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	414	MY8	C14-C13-C11	-2.55	110.32	116.22
3	B	414	MY8	O1-C29-C28	-2.53	105.89	111.39
2	A	413	HEM	CHB-C4A-NA	2.50	128.40	123.86
2	A	413	HEM	CHA-C4D-C3D	-2.49	120.63	125.23
2	B	413	HEM	CHB-C1B-NB	2.49	127.45	124.37
3	B	414	MY8	O3-C16-C25	-2.48	105.94	110.93
2	A	413	HEM	CMD-C2D-C1D	2.47	128.90	125.03
3	B	414	MY8	C14-C13-C11	-2.47	110.52	116.22
5	A	1391	SPD	C7-N6-C5	2.45	125.02	113.40
2	A	413	HEM	CMC-C2C-C3C	-2.39	122.64	128.43
2	A	413	HEM	CMB-C2B-C1B	2.35	128.70	125.03
3	A	414	MY8	O5-C24-C21	-2.33	105.59	109.87
3	B	414	MY8	C17-O3-C16	-2.33	112.46	117.98
2	B	413	HEM	C1A-CHA-C4D	-2.31	120.81	126.25
3	B	414	MY8	O5-C24-C21	-2.30	105.65	109.87
2	A	413	HEM	O2D-CGD-CBD	2.26	121.14	114.00
3	B	414	MY8	O5-C24-C17	-2.25	104.72	110.08
2	A	413	HEM	CHD-C4C-C3C	2.20	128.92	125.21
2	B	413	HEM	CAD-CBD-CGD	-2.15	107.95	113.67
3	A	414	MY8	C20-C21-C24	-2.15	106.93	110.02
2	B	413	HEM	CAA-C2A-C1A	2.10	129.04	124.94
3	A	414	MY8	O6-C29-C28	2.05	130.66	123.69
2	B	413	HEM	CMC-C2C-C1C	2.04	128.32	124.73

There are no chirality outliers.

All (31) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	414	MY8	O1-C1-C4-C6
3	A	414	MY8	C11-C10-C9-C8
3	A	414	MY8	C9-C10-C11-C13
3	A	414	MY8	O2-C10-C11-C12
3	A	414	MY8	O2-C10-C11-C13
3	B	414	MY8	C11-C10-C9-C8
3	B	414	MY8	O2-C10-C11-C12
3	B	414	MY8	O2-C10-C11-C13
4	A	1390	GOL	C1-C2-C3-O3
5	A	1391	SPD	N6-C7-C8-C9
5	A	1391	SPD	C4-C5-N6-C7
3	A	414	MY8	C5-C4-C6-C7
3	B	414	MY8	C5-C4-C6-C7
4	A	1390	GOL	O2-C2-C3-O3

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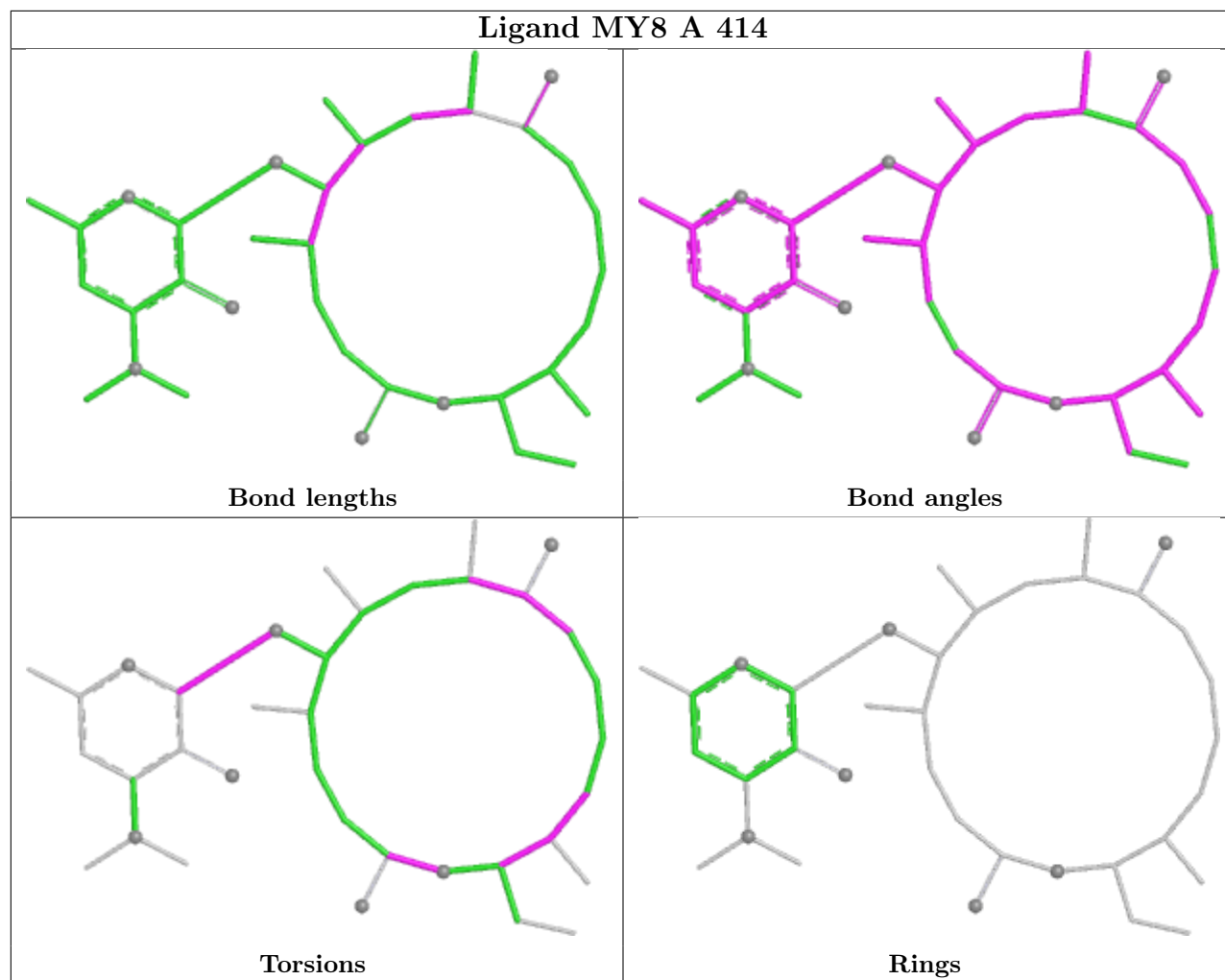
Mol	Chain	Res	Type	Atoms
5	A	1392	SPD	N6-C7-C8-C9
5	A	1392	SPD	C2-C3-C4-C5
5	A	1392	SPD	C3-C4-C5-N6
5	A	1391	SPD	C2-C3-C4-C5
3	A	414	MY8	O6-C29-O1-C1
2	B	413	HEM	C4C-C3C-CAC-CBC
5	A	1392	SPD	C8-C7-N6-C5
3	B	414	MY8	C20-C21-N1-C23
3	A	414	MY8	C24-C17-O3-C16
5	A	1393	SPD	C8-C7-N6-C5
3	B	414	MY8	C20-C21-N1-C22
5	A	1391	SPD	C8-C7-N6-C5
3	A	414	MY8	C28-C29-O1-C1
3	B	414	MY8	C2-C1-C4-C5
5	A	1392	SPD	C4-C5-N6-C7
3	B	414	MY8	C24-C17-O3-C16
3	B	414	MY8	C9-C10-C11-C12

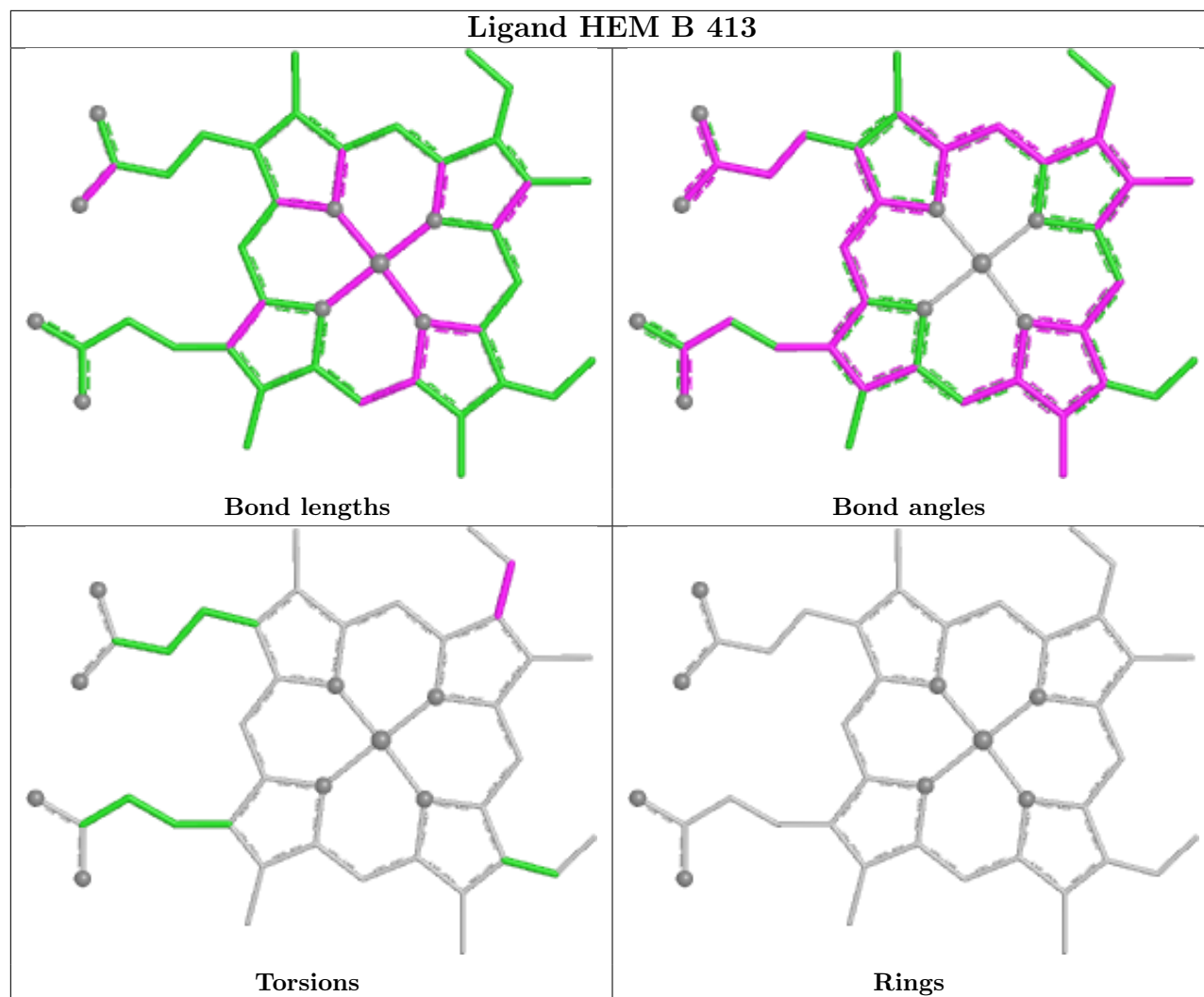
There are no ring outliers.

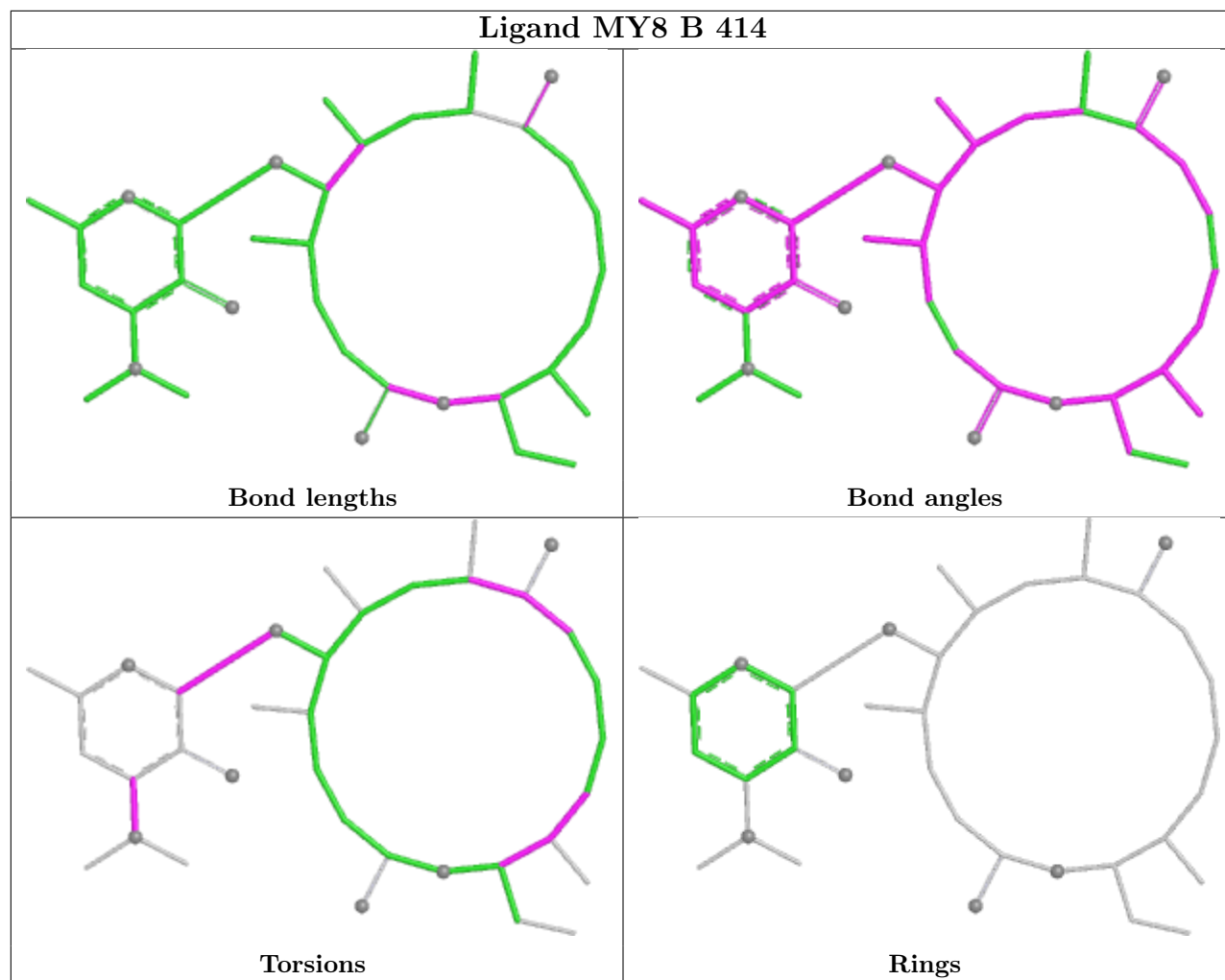
6 monomers are involved in 17 short contacts:

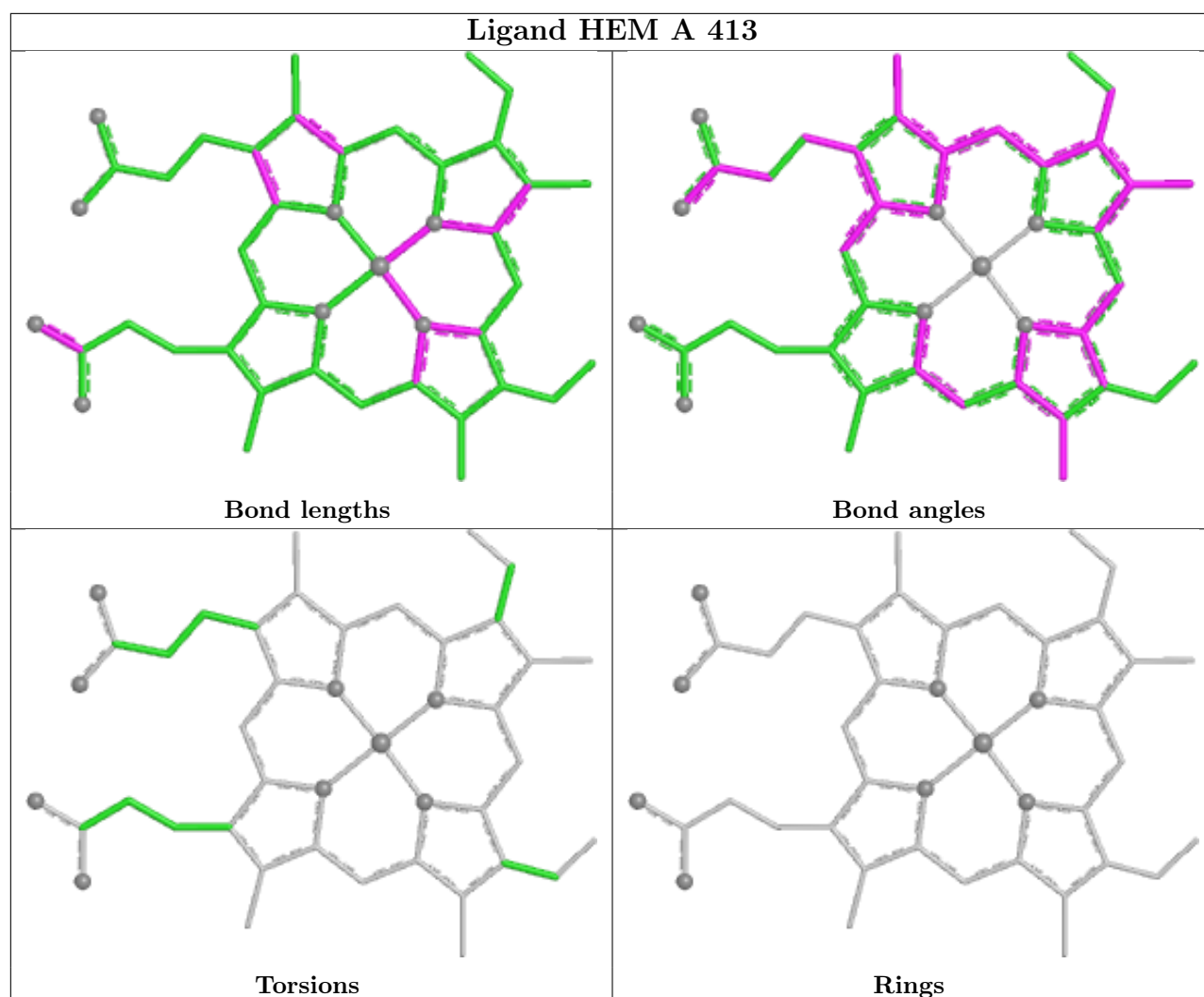
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	1392	SPD	3	0
3	A	414	MY8	1	0
5	A	1391	SPD	3	0
4	A	1390	GOL	3	0
2	B	413	HEM	8	0
2	A	413	HEM	2	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.









5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

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

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	389/408 (95%)	-0.44	0 100 100	15, 26, 44, 59	0
1	B	388/408 (95%)	-0.46	2 (0%) 87 86	14, 26, 43, 61	2 (0%)
All	All	777/816 (95%)	-0.45	2 (0%) 90 89	14, 26, 44, 61	2 (0%)

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	26	THR	2.8
1	B	25	ALA	2.1

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [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, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q < 0.9' lists the number of atoms with occupancy less than 0.9.

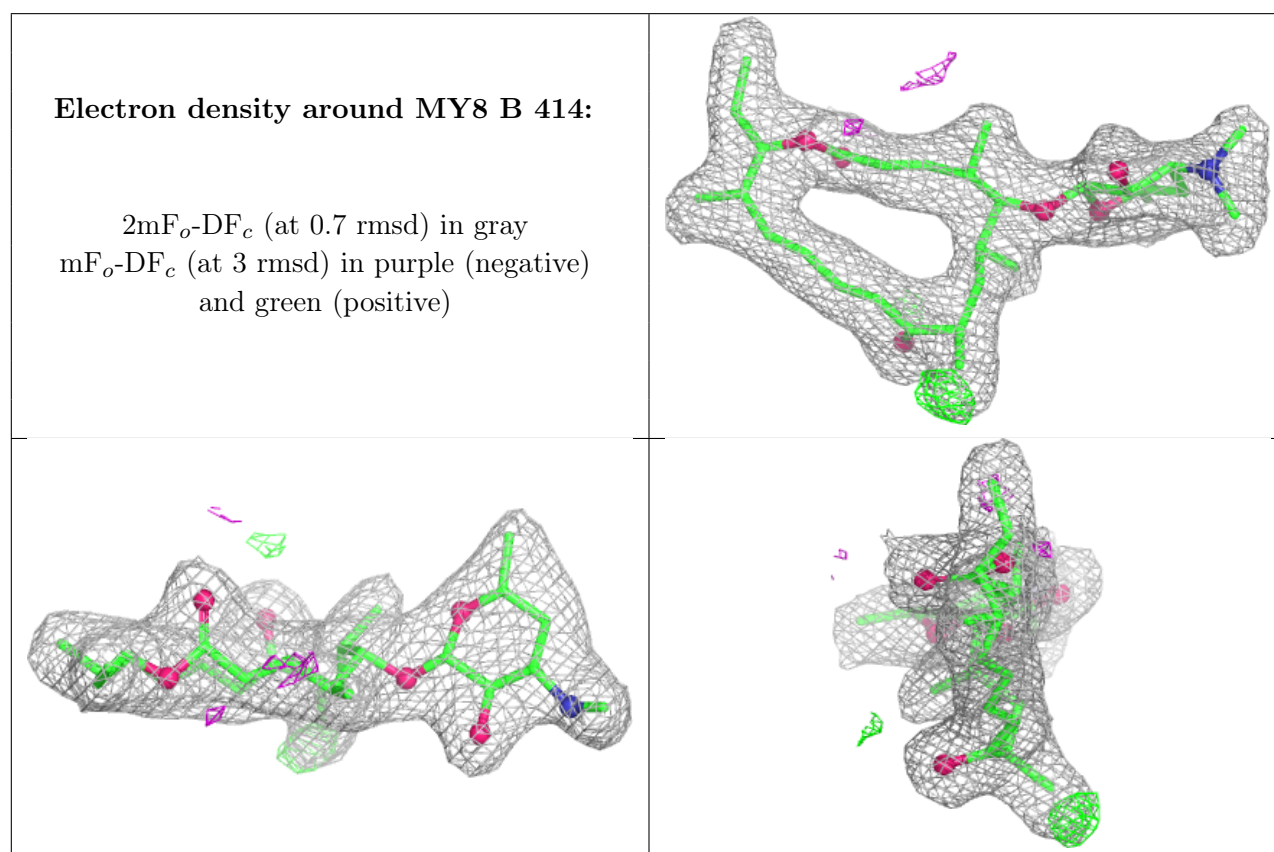
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	SPD	A	1393	10/10	0.81	0.13	48,55,56,59	0
4	GOL	A	1390	6/6	0.82	0.12	45,50,52,53	0
5	SPD	A	1392	10/10	0.83	0.15	39,42,47,47	0

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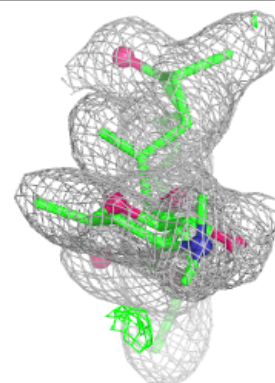
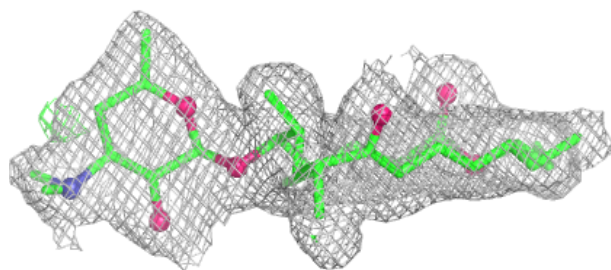
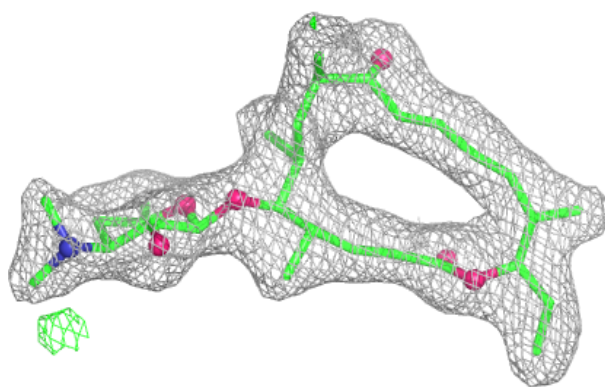
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	SPD	A	1391	10/10	0.88	0.14	29,31,34,34	0
4	GOL	A	1389	6/6	0.95	0.07	32,37,40,42	0
3	MY8	B	414	36/36	0.96	0.05	18,21,24,24	0
3	MY8	A	414	36/36	0.96	0.06	16,19,24,25	0
2	HEM	B	413	43/43	0.98	0.06	12,17,19,23	0
2	HEM	A	413	43/43	0.98	0.05	12,16,18,19	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.



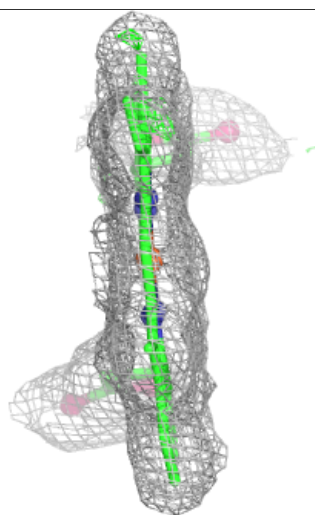
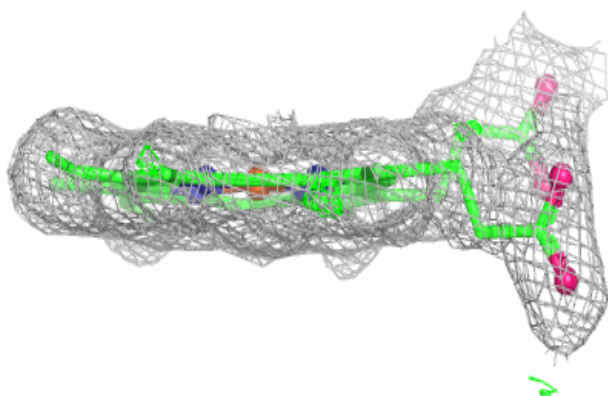
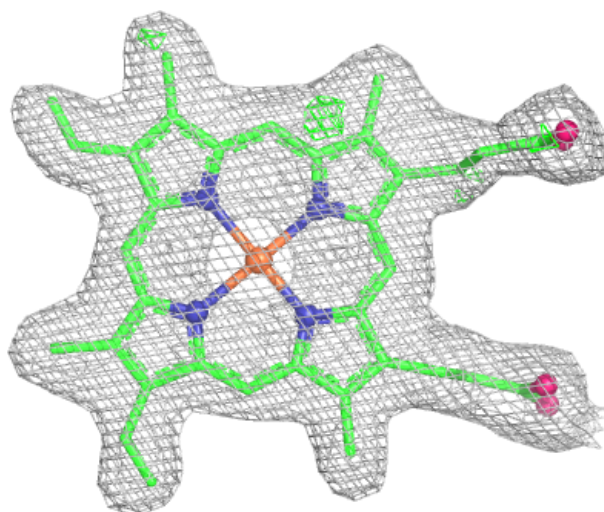
Electron density around MY8 A 414:

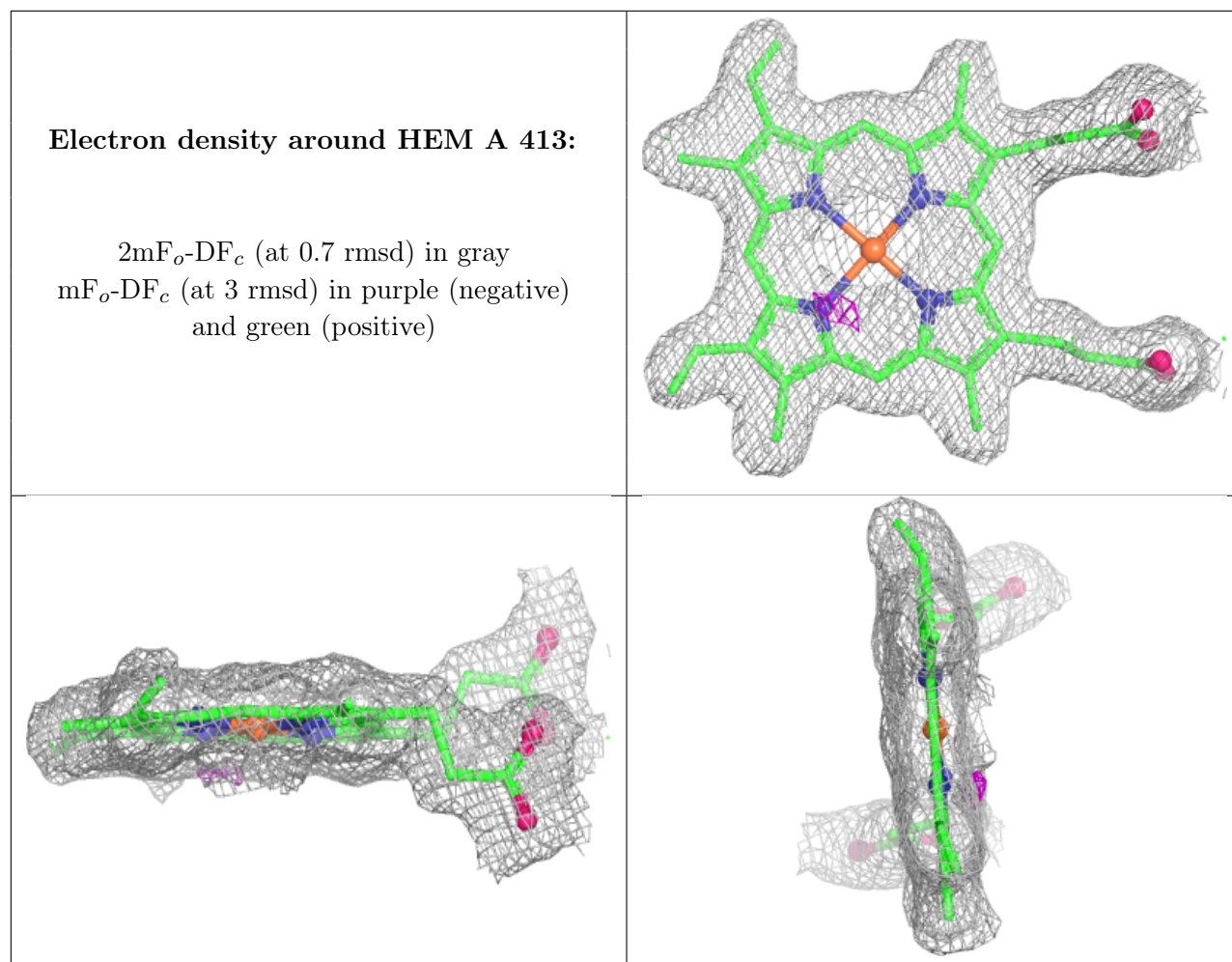
$2mF_o - DF_c$ (at 0.7 rmsd) in gray
 $mF_o - DF_c$ (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around HEM B 413:

$2mF_o-DF_c$ (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.