



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 02:32 AM UTC

PDB ID : 4PCU / pdb_00004pcu
Title : Crystal structure of delta516-525 E201S human cystathionine beta-synthase with AdoMet
Authors : Ereno-Orbea, J.; Majtan, T.; Oyenarte, I.; Kraus, J.P.; Martinez-Cruz, L.A.
Deposited on : 2014-04-16
Resolution : 3.58 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)
Xtrriage (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

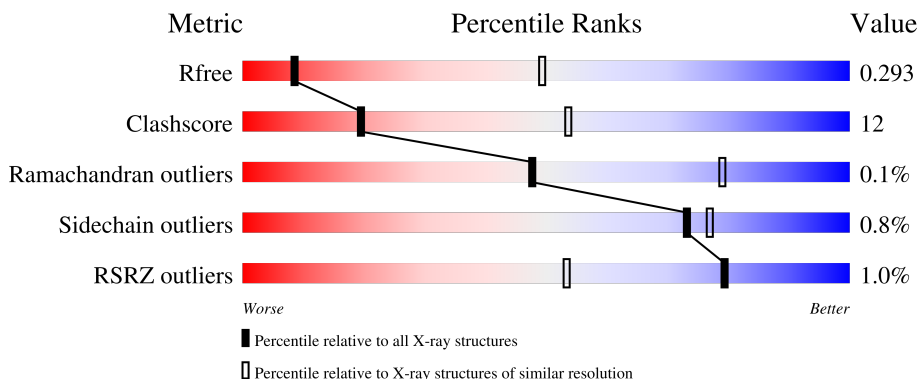
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION



The reported resolution of this entry is 3.58 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	1521 (3.66-3.50)
Clashscore	190562	1595 (3.66-3.50)
Ramachandran outliers	187476	1551 (3.66-3.50)
Sidechain outliers	187428	1551 (3.66-3.50)
RSRZ outliers	180081	1520 (3.66-3.50)

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	549	 70% 17% 11%
1	B	549	 74% 14% 11%

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 7565 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 Cystathionine beta-synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	486	3681	2335	638	687	21	0	0	0
1	B	489	3714	2352	644	697	21	0	0	0

There are 40 discrepancies between the modelled and reference sequences:

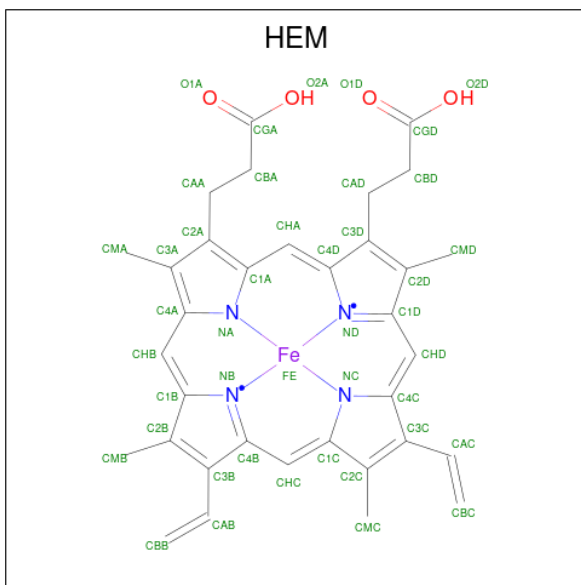
Chain	Residue	Modelled	Actual	Comment	Reference
A	2	GLY	PRO	conflict	UNP P35520
A	201	SER	GLU	engineered mutation	UNP P35520
A	?	-	ILE	deletion	UNP P35520
A	?	-	GLN	deletion	UNP P35520
A	?	-	TYR	deletion	UNP P35520
A	?	-	HIS	deletion	UNP P35520
A	?	-	SER	deletion	UNP P35520
A	?	-	THR	deletion	UNP P35520
A	?	-	GLY	deletion	UNP P35520
A	?	-	LYS	deletion	UNP P35520
A	?	-	SER	deletion	UNP P35520
A	?	-	SER	deletion	UNP P35520
A	552	GLU	-	expression tag	UNP P35520
A	553	LEU	-	expression tag	UNP P35520
A	554	HIS	-	expression tag	UNP P35520
A	555	HIS	-	expression tag	UNP P35520
A	556	HIS	-	expression tag	UNP P35520
A	557	HIS	-	expression tag	UNP P35520
A	558	HIS	-	expression tag	UNP P35520
A	559	HIS	-	expression tag	UNP P35520
B	2	GLY	PRO	conflict	UNP P35520
B	201	SER	GLU	engineered mutation	UNP P35520
B	?	-	ILE	deletion	UNP P35520
B	?	-	GLN	deletion	UNP P35520
B	?	-	TYR	deletion	UNP P35520

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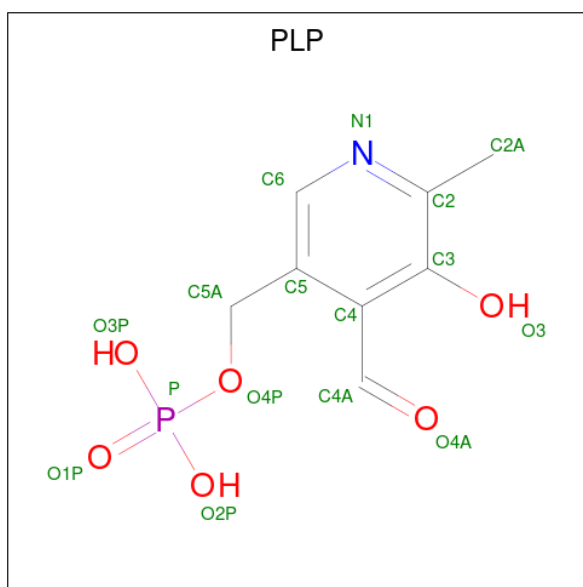
Chain	Residue	Modelled	Actual	Comment	Reference
B	?	-	HIS	deletion	UNP P35520
B	?	-	SER	deletion	UNP P35520
B	?	-	THR	deletion	UNP P35520
B	?	-	GLY	deletion	UNP P35520
B	?	-	LYS	deletion	UNP P35520
B	?	-	SER	deletion	UNP P35520
B	?	-	SER	deletion	UNP P35520
B	552	GLU	-	expression tag	UNP P35520
B	553	LEU	-	expression tag	UNP P35520
B	554	HIS	-	expression tag	UNP P35520
B	555	HIS	-	expression tag	UNP P35520
B	556	HIS	-	expression tag	UNP P35520
B	557	HIS	-	expression tag	UNP P35520
B	558	HIS	-	expression tag	UNP P35520
B	559	HIS	-	expression tag	UNP P35520

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



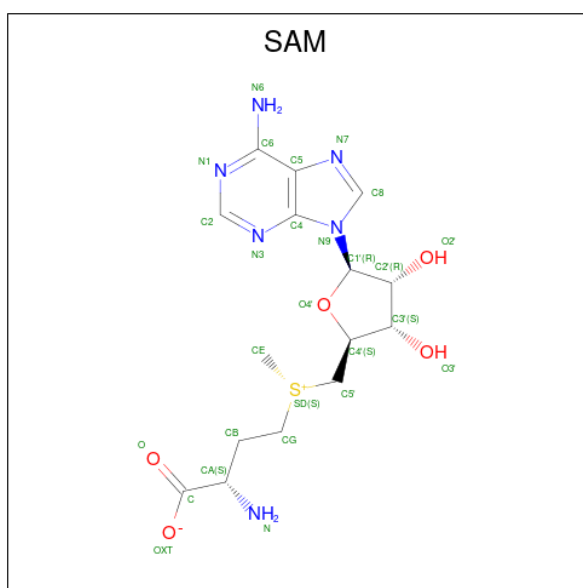
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
2	A	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
2	B	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		

- Molecule 3 is PYRIDOXAL-5'-PHOSPHATE (CCD ID: PLP) (formula: $C_8H_{10}NO_6P$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	
			Total	C	N	O	P			
3	A	1	Total	15	8	1	5	1	0	0
3	B	1	Total	15	8	1	5	1	0	0

- Molecule 4 is S-ADENOSYLMETHIONINE (CCD ID: SAM) (formula: $C_{15}H_{22}N_6O_5S$).



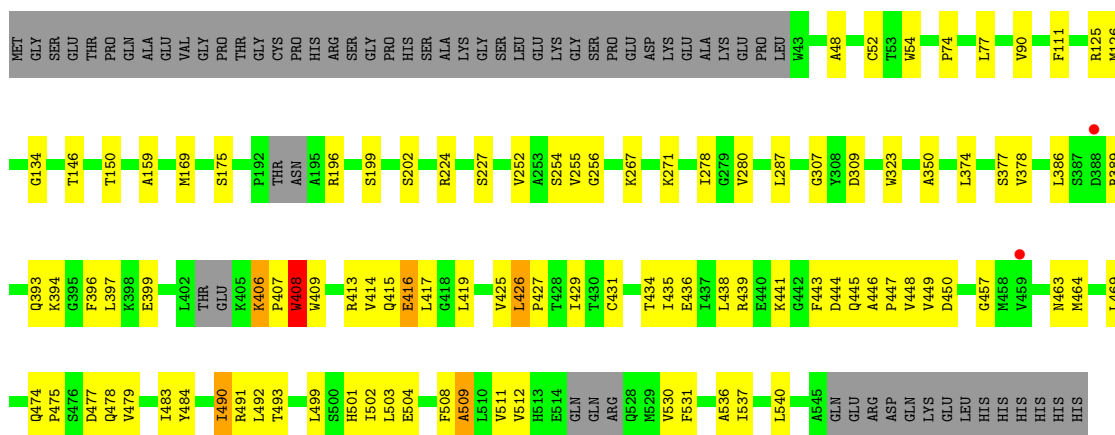
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	
			Total	C	N	O	S			
4	A	1	Total	27	15	6	5	1	0	0
4	B	1	Total	27	15	6	5	1	0	0

3 Residue-property plots


These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

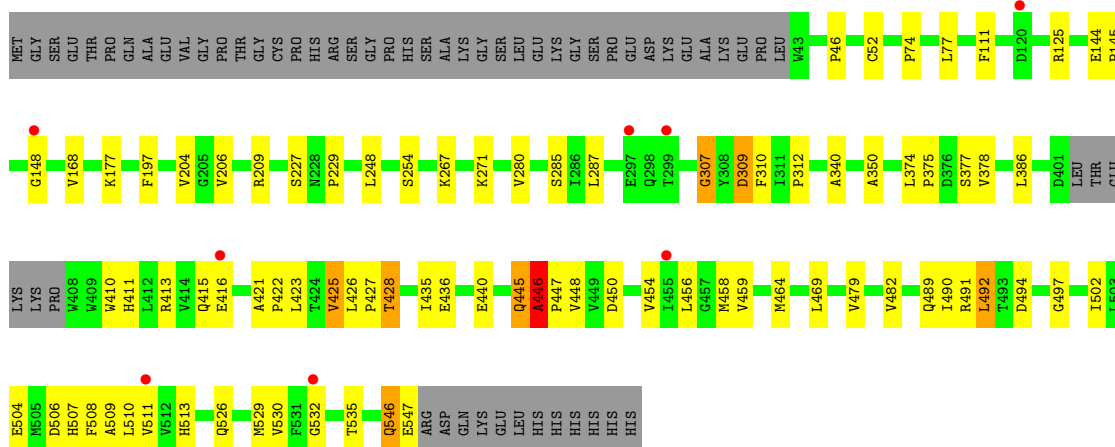
- Molecule 1: Cystathionine beta-synthase

Chain A: 



- Molecule 1: Cystathionine beta-synthase

Chain B: 



4 Data and refinement statistics i

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	141.35Å 141.35Å 108.53Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	40.80 – 3.58 40.80 – 3.58	Depositor EDS
% Data completeness (in resolution range)	98.2 (40.80-3.58) 98.4 (40.80-3.58)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.21 (at 3.57Å)	Xtrriage
Refinement program	PHENIX (phenix.refine: 1.9_1692)	Depositor
R, R_{free}	0.257 , 0.277 (Not available) , 0.293	Depositor DCC
R_{free} test set	756 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	105.5	Xtrriage
Anisotropy	0.456	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 86.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.27$	Xtrriage
Estimated twinning fraction	0.057 for -h,-k,l	Xtrriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	7565	wwPDB-VP
Average B, all atoms (Å ²)	117.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.74% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality i

5.1 Standard geometry i

Bond lengths and bond angles in the following residue types are not validated in this section: HEM, SAM, PLP

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	0.50	1/3751 (0.0%)	0.90	17/5091 (0.3%)
1	B	0.44	0/3784	0.99	13/5135 (0.3%)
All	All	0.47	1/7535 (0.0%)	0.95	30/10226 (0.3%)

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	A	0	1
1	B	0	2
All	All	0	3

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	408	TRP	C-N	17.11	1.59	1.33

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	428	THR	O-C-N	-21.92	92.21	122.46
1	A	399	GLU	N-CA-C	15.94	137.95	107.98
1	B	445	GLN	N-CA-C	15.87	129.19	108.24
1	B	428	THR	CA-C-N	13.61	140.42	120.95
1	B	428	THR	C-N-CA	13.61	140.42	120.95
1	B	307	GLY	N-CA-C	9.24	127.58	114.64
1	B	446	ALA	CA-C-N	-8.83	110.63	119.90
1	B	446	ALA	C-N-CA	-8.83	110.63	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	546	GLN	O-C-N	-6.90	114.06	122.34
1	A	408	TRP	N-CA-C	6.64	120.97	112.34
1	A	509	ALA	CA-C-N	-6.58	112.89	122.44
1	A	509	ALA	C-N-CA	-6.58	112.89	122.44
1	A	490	ILE	CA-C-N	6.22	134.38	121.32
1	A	490	ILE	C-N-CA	6.22	134.38	121.32
1	A	408	TRP	CA-C-N	-5.88	110.19	121.94
1	A	408	TRP	C-N-CA	-5.88	110.19	121.94
1	B	421	ALA	CA-C-N	5.87	125.88	119.89
1	B	421	ALA	C-N-CA	5.87	125.88	119.89
1	A	457	GLY	N-CA-C	5.81	120.19	111.42
1	B	445	GLN	CA-C-O	5.77	127.11	121.05
1	A	508	PHE	N-CA-C	5.75	116.14	108.38
1	A	175	SER	N-CA-C	5.51	118.81	111.75
1	A	399	GLU	CB-CA-C	-5.45	102.53	111.68
1	A	169	MET	N-CA-C	5.43	112.91	108.07
1	A	408	TRP	O-C-N	5.38	129.32	122.33
1	A	134	GLY	N-CA-C	-5.36	107.92	115.32
1	B	492	LEU	N-CA-C	5.22	117.37	111.11
1	B	309	ASP	N-CA-C	5.14	121.75	110.80
1	A	399	GLU	N-CA-CB	-5.13	102.90	110.49
1	A	406	LYS	N-CA-C	5.01	112.53	108.07

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	408	TRP	Mainchain
1	B	446	ALA	Peptide
1	B	546	GLN	Mainchain

5.2 Too-close contacts [\(i\)](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3681	0	3659	102	1
1	B	3714	0	3701	82	3

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	43	0	30	4	0
2	B	43	0	30	4	0
3	A	15	0	7	1	0
3	B	15	0	6	0	0
4	A	27	0	22	4	0
4	B	27	0	22	6	0
All	All	7565	0	7477	174	3

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 12.

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:A:435:ILE:O	1:A:439:ARG:HG3	1.22	1.32
1:A:425:VAL:HG22	1:A:447:PRO:O	1.34	1.25
1:B:513:HIS:HD2	1:B:529:MET:HB3	1.06	1.14
1:A:415:GLN:HB3	1:A:492:LEU:CD1	1.80	1.09
1:A:414:VAL:HG13	1:A:417:LEU:HD12	1.20	1.08
1:B:513:HIS:CD2	1:B:529:MET:HB3	1.94	1.02
1:A:415:GLN:HB3	1:A:492:LEU:HD12	1.40	1.01
1:B:513:HIS:HD2	1:B:529:MET:CB	1.76	0.97
1:B:445:GLN:NE2	4:B:603:SAM:N7	2.21	0.88
1:B:427:PRO:CD	1:B:428:THR:H	1.87	0.87
1:A:504:GLU:OE2	1:B:469:LEU:HD12	1.75	0.85
1:A:540:LEU:HB3	1:B:435:ILE:HD13	1.60	0.83
1:B:427:PRO:HD2	1:B:428:THR:H	1.42	0.82
1:A:537:ILE:CD1	1:B:435:ILE:HG23	2.10	0.82
1:A:426:LEU:O	1:A:429:ILE:HG12	1.80	0.82
1:A:435:ILE:O	1:A:439:ARG:CG	2.19	0.81
1:A:504:GLU:OE2	1:B:469:LEU:CD1	2.29	0.81
1:A:146:THR:HG21	1:A:150:THR:HB	1.62	0.79
1:A:425:VAL:CG2	1:A:447:PRO:O	2.27	0.78
1:A:406:LYS:CB	1:A:407:PRO:HD2	2.15	0.77
1:A:415:GLN:HB3	1:A:492:LEU:HD11	1.63	0.76
1:A:491:ARG:NH1	1:A:493:THR:OG1	2.20	0.74
1:A:414:VAL:HG13	1:A:417:LEU:CD1	2.10	0.73
1:A:415:GLN:CB	1:A:492:LEU:CD1	2.64	0.72
1:B:425:VAL:HG13	1:B:446:ALA:HB2	1.72	0.72
1:A:436:GLU:HA	1:A:439:ARG:HD3	1.72	0.71
1:A:434:THR:HG22	1:A:438:LEU:HD12	1.71	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:537:ILE:HD11	1:B:435:ILE:HG23	1.74	0.70
1:A:406:LYS:CB	1:A:407:PRO:CD	2.70	0.70
1:A:393:GLN:OE1	1:A:501:HIS:O	2.10	0.69
1:A:415:GLN:CB	1:A:492:LEU:HD12	2.19	0.69
1:B:427:PRO:CD	1:B:428:THR:N	2.57	0.66
1:B:491:ARG:HA	1:B:511:VAL:HA	1.77	0.66
1:B:513:HIS:O	1:B:526:GLN:N	2.29	0.66
1:B:410:TRP:HD1	1:B:411:HIS:HD2	1.43	0.66
1:B:490:ILE:HB	1:B:494:ASP:OD2	1.96	0.66
1:A:393:GLN:NE2	1:A:504:GLU:OE1	2.22	0.64
1:A:425:VAL:HG23	1:A:448:VAL:HA	1.79	0.64
2:B:601:HEM:HMB1	2:B:601:HEM:HBB2	1.80	0.63
1:A:502:ILE:HG21	1:A:509:ALA:HB2	1.80	0.62
1:B:145:PRO:HB3	1:B:204:VAL:HA	1.82	0.62
1:B:285:SER:OG	1:B:307:GLY:O	2.16	0.61
1:A:537:ILE:HD12	1:B:435:ILE:HG23	1.79	0.61
1:B:447:PRO:HA	1:B:458:MET:HG2	1.82	0.60
1:A:389:ARG:O	1:A:393:GLN:HG2	2.01	0.60
1:A:415:GLN:CB	1:A:492:LEU:HD11	2.27	0.59
1:B:426:LEU:HB3	1:B:427:PRO:HD2	1.84	0.59
1:B:427:PRO:CG	1:B:428:THR:N	2.66	0.59
1:A:474:GLN:N	1:A:477:ASP:OD2	2.36	0.58
1:A:512:VAL:HG12	1:A:530:VAL:HA	1.84	0.58
1:B:450:ASP:HB3	1:B:456:LEU:HD21	1.85	0.58
1:B:513:HIS:CD2	1:B:529:MET:CB	2.69	0.57
1:A:254:SER:HA	1:A:280:VAL:HB	1.86	0.57
1:A:436:GLU:HA	1:A:439:ARG:CD	2.34	0.57
1:A:436:GLU:O	1:A:439:ARG:HB2	2.04	0.57
2:A:601:HEM:HMB1	2:A:601:HEM:HBB2	1.87	0.57
1:B:490:ILE:HG12	1:B:509:ALA:HB1	1.85	0.57
1:B:427:PRO:CG	1:B:428:THR:H	2.18	0.57
1:A:408:TRP:CZ3	1:A:409:TRP:HB3	2.39	0.56
1:B:148:GLY:HA3	1:B:177:LYS:HD3	1.86	0.56
1:B:254:SER:HA	1:B:280:VAL:HB	1.87	0.56
1:B:490:ILE:HD13	1:B:502:ILE:HD12	1.88	0.56
2:B:601:HEM:HMC1	2:B:601:HEM:HBC2	1.88	0.55
1:A:445:GLN:NE2	4:A:603:SAM:O	2.31	0.55
4:A:603:SAM:H5'2	4:A:603:SAM:OXT	2.07	0.55
1:A:540:LEU:HD13	1:B:435:ILE:HD13	1.89	0.55
1:A:474:GLN:HG3	1:A:477:ASP:OD2	2.07	0.54
1:A:429:ILE:CG1	1:A:479:VAL:HG11	2.37	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:435:ILE:HD11	1:B:464:MET:HG2	1.91	0.53
1:A:438:LEU:HD21	1:A:446:ALA:HB2	1.90	0.53
1:A:540:LEU:CB	1:B:435:ILE:HD13	2.37	0.52
1:B:507:HIS:HB2	1:B:508:PHE:HD2	1.74	0.52
1:A:74:PRO:HG2	1:A:77:LEU:HD23	1.90	0.52
1:A:52:CYS:HA	2:A:601:HEM:C1A	2.45	0.52
1:A:413:ARG:HD3	1:A:493:THR:O	2.10	0.52
1:B:267:LYS:HD2	1:B:271:LYS:HE2	1.92	0.52
1:B:427:PRO:HG2	1:B:428:THR:HG23	1.91	0.52
1:A:393:GLN:OE1	1:A:504:GLU:HB2	2.11	0.51
1:A:350:ALA:HB1	1:A:374:LEU:HD22	1.93	0.51
1:B:74:PRO:HG2	1:B:77:LEU:HD23	1.92	0.50
1:B:125:ARG:HG3	1:B:227:SER:HB3	1.92	0.50
1:A:287:LEU:HD12	1:A:307:GLY:HA2	1.92	0.50
1:B:410:TRP:CD1	1:B:411:HIS:HD2	2.25	0.50
1:B:448:VAL:O	1:B:456:LEU:N	2.39	0.50
1:B:436:GLU:O	1:B:440:GLU:HG3	2.12	0.50
1:A:414:VAL:HG12	1:A:414:VAL:O	2.11	0.50
2:A:601:HEM:HBC2	2:A:601:HEM:HMC1	1.91	0.50
1:A:267:LYS:HD2	1:A:271:LYS:HE2	1.93	0.50
1:B:435:ILE:HG22	1:B:435:ILE:O	2.12	0.50
1:B:510:LEU:HD11	1:B:530:VAL:HG12	1.93	0.50
1:B:350:ALA:HB1	1:B:374:LEU:HD22	1.93	0.49
1:B:502:ILE:O	1:B:506:ASP:N	2.33	0.49
1:B:410:TRP:CE2	1:B:497:GLY:HA3	2.47	0.49
1:A:475:PRO:HG2	1:B:547:GLU:HB2	1.95	0.49
1:A:414:VAL:HG22	1:A:499:LEU:HD22	1.94	0.49
1:B:489:GLN:O	1:B:489:GLN:HG3	2.13	0.49
1:A:394:LYS:HB3	1:A:396:PHE:CE2	2.48	0.49
4:B:603:SAM:H8	4:B:603:SAM:O	2.13	0.48
1:A:427:PRO:HD3	1:A:449:VAL:O	2.14	0.48
1:A:394:LYS:CB	1:A:396:PHE:CE2	2.97	0.48
1:B:46:PRO:HG3	1:B:197:PHE:HZ	1.77	0.48
1:A:490:ILE:HD13	1:A:499:LEU:HB2	1.94	0.48
1:A:530:VAL:HG12	1:A:531:PHE:N	2.29	0.48
1:A:490:ILE:HG21	1:A:499:LEU:HD13	1.96	0.47
1:B:413:ARG:HB2	1:B:415:GLN:HG2	1.95	0.47
1:B:535:THR:HG21	4:B:603:SAM:O	2.15	0.47
1:A:511:VAL:O	1:A:531:PHE:N	2.46	0.47
1:A:443:PHE:HD1	4:A:603:SAM:SD	2.38	0.46
1:A:463:ASN:HD22	1:A:484:TYR:HB2	1.79	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:540:LEU:HD13	1:B:435:ILE:CD1	2.45	0.46
4:B:603:SAM:O	4:B:603:SAM:H5'2	2.16	0.46
1:A:431:CYS:HB3	1:A:464:MET:HE1	1.96	0.46
1:A:537:ILE:HD12	1:B:435:ILE:CG2	2.46	0.46
1:B:423:LEU:HD23	1:B:423:LEU:C	2.40	0.46
1:B:479:VAL:O	1:B:482:VAL:HG22	2.16	0.46
1:B:52:CYS:HA	2:B:601:HEM:C1A	2.51	0.46
1:A:448:VAL:HG11	1:A:483:ILE:HG12	1.98	0.45
1:A:378:VAL:HG21	1:A:386:LEU:HD13	1.99	0.45
1:A:503:LEU:HD13	1:A:536:ALA:HA	1.99	0.45
1:B:446:ALA:HB3	1:B:459:VAL:HG12	1.98	0.45
1:B:450:ASP:OD2	1:B:454:VAL:HB	2.17	0.45
1:A:125:ARG:HG3	1:A:227:SER:HB3	1.98	0.45
1:A:425:VAL:CG2	1:A:448:VAL:HA	2.44	0.45
1:A:504:GLU:OE2	1:B:469:LEU:HD13	2.14	0.45
1:A:537:ILE:CD1	1:B:435:ILE:CG2	2.90	0.45
1:B:111:PHE:HB2	1:B:377:SER:HB3	1.99	0.45
1:A:469:LEU:HD12	1:B:504:GLU:OE2	2.17	0.45
1:A:444:ASP:CG	4:B:603:SAM:HN2	2.23	0.45
1:A:490:ILE:HG22	1:A:509:ALA:CB	2.47	0.45
1:A:196:ARG:HB3	1:A:199:SER:HB2	1.97	0.44
1:A:427:PRO:HD3	1:A:450:ASP:HA	1.99	0.44
1:A:426:LEU:HB2	1:A:429:ILE:HG23	1.99	0.44
1:A:255:VAL:HG21	1:A:323:TRP:CZ3	2.53	0.44
1:B:425:VAL:CG1	1:B:446:ALA:HB2	2.45	0.44
4:A:603:SAM:OXT	4:A:603:SAM:H8	2.16	0.44
1:B:427:PRO:HG2	1:B:428:THR:N	2.32	0.44
1:B:310:PHE:O	1:B:312:PRO:HD3	2.18	0.44
1:B:378:VAL:HG21	1:B:386:LEU:HD13	1.99	0.44
1:A:434:THR:HG22	1:A:438:LEU:CD1	2.43	0.43
1:A:540:LEU:HB3	1:B:435:ILE:HG21	1.99	0.43
1:A:111:PHE:HB2	1:A:377:SER:HB3	2.00	0.43
1:A:126:MET:HG2	1:A:227:SER:HB2	2.00	0.43
1:A:464:MET:HE2	1:A:464:MET:HB3	1.71	0.43
1:B:445:GLN:HG2	4:B:603:SAM:HN61	1.84	0.43
1:B:168:VAL:HG11	1:B:206:VAL:HG13	2.00	0.43
1:A:256:GLY:HA3	3:A:602:PLP:H5A1	2.01	0.43
1:B:374:LEU:HA	1:B:375:PRO:HD3	1.88	0.43
1:A:419:LEU:HD12	1:A:511:VAL:HG21	2.00	0.42
1:B:287:LEU:HD12	1:B:307:GLY:HA2	2.01	0.42
1:A:309:ASP:OD1	1:A:309:ASP:N	2.52	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:425:VAL:HB	1:A:429:ILE:CD1	2.49	0.42
1:A:491:ARG:HH11	1:A:491:ARG:HG2	1.84	0.42
1:B:422:PRO:HD3	1:B:532:GLY:HA2	2.01	0.42
1:B:492:LEU:N	1:B:511:VAL:HG13	2.33	0.42
1:B:229:PRO:HB2	2:B:601:HEM:HBC1	2.02	0.42
1:A:48:ALA:HB3	1:A:224:ARG:NH1	2.35	0.42
1:A:90:VAL:HG22	1:B:77:LEU:HD12	2.01	0.41
1:A:441:LYS:HB3	1:A:441:LYS:HE3	1.86	0.41
1:B:168:VAL:HG11	1:B:206:VAL:CG1	2.50	0.41
1:B:464:MET:HE1	1:B:479:VAL:HG23	2.01	0.41
1:A:159:ALA:HB1	1:B:340:ALA:O	2.20	0.41
1:A:414:VAL:HA	1:A:417:LEU:HG	2.03	0.41
1:A:425:VAL:HB	1:A:429:ILE:HD11	2.01	0.41
1:A:429:ILE:HG12	1:A:479:VAL:HG11	2.03	0.41
1:A:408:TRP:CD1	1:A:408:TRP:H	2.39	0.41
1:A:463:ASN:ND2	1:A:484:TYR:HB2	2.36	0.41
1:A:252:VAL:HG22	1:A:278:ILE:HB	2.03	0.41
1:A:414:VAL:O	1:A:414:VAL:CG1	2.70	0.40
1:B:46:PRO:HG3	1:B:197:PHE:CZ	2.56	0.40
1:B:144:GLU:HA	1:B:145:PRO:HD3	1.80	0.40
1:B:426:LEU:HB3	1:B:427:PRO:CD	2.50	0.40
1:A:54:TRP:HB2	2:A:601:HEM:C4B	2.55	0.40
1:A:199:SER:HB3	1:A:202:SER:HB3	2.03	0.40
1:A:540:LEU:HD22	1:B:435:ILE:CD1	2.51	0.40

All (3) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:416:GLU:OE1	1:B:416:GLU:OE1[6_555]	1.60	0.60
1:A:416:GLU:OE2	1:B:209:ARG:NE[5_554]	1.93	0.27
1:B:416:GLU:CD	1:B:416:GLU:OE1[6_555]	2.18	0.02

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	A	478/549 (87%)	460 (96%)	18 (4%)	0	100	100
1	B	483/549 (88%)	467 (97%)	15 (3%)	1 (0%)	43	72
All	All	961/1098 (88%)	927 (96%)	33 (3%)	1 (0%)	48	79

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	309	ASP

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	389/463 (84%)	385 (99%)	4 (1%)	68	75
1	B	396/463 (86%)	394 (100%)	2 (0%)	81	80
All	All	785/926 (85%)	779 (99%)	6 (1%)	73	77

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	397	LEU
1	A	416	GLU
1	A	426	LEU
1	A	478	GLN
1	B	248	LEU
1	B	425	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	380	ASN

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Mol	Chain	Res	Type
1	B	149	ASN
1	B	380	ASN
1	B	411	HIS
1	B	445	GLN
1	B	474	GLN
1	B	513	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

6 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	A	601	1	50,50,50	1.62	10 (20%)	67,82,82	1.20	8 (11%)
4	SAM	B	603	-	27,29,29	1.33	2 (7%)	34,42,42	2.45	12 (35%)
3	PLP	A	602	1	15,15,16	0.95	1 (6%)	21,22,23	1.41	4 (19%)
4	SAM	A	603	-	27,29,29	1.27	5 (18%)	34,42,42	1.83	12 (35%)
3	PLP	B	602	-	15,15,16	1.00	1 (6%)	21,22,23	1.45	3 (14%)
2	HEM	B	601	1	50,50,50	1.65	10 (20%)	67,82,82	1.19	6 (8%)

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	A	601	1	-	0/14/54/54	-
4	SAM	B	603	-	-	5/17/33/33	0/3/3/3
3	PLP	A	602	1	-	0/6/6/8	0/1/1/1
4	SAM	A	603	-	-	6/17/33/33	0/3/3/3
3	PLP	B	602	-	-	0/6/6/8	0/1/1/1
2	HEM	B	601	1	-	0/14/54/54	-

All (29) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	601	HEM	C3D-C2D	6.98	1.51	1.36
2	A	601	HEM	C3D-C2D	6.89	1.51	1.36
4	B	603	SAM	C5-C4	3.56	1.45	1.39
4	B	603	SAM	C4-N9	-3.41	1.30	1.37
2	A	601	HEM	CAB-C3B	3.15	1.55	1.47
2	B	601	HEM	FE-ND	3.08	2.04	1.94
2	B	601	HEM	FE-NA	3.04	2.05	1.95
2	B	601	HEM	CAC-C3C	2.93	1.55	1.47
2	B	601	HEM	CAB-C3B	2.91	1.55	1.47
2	A	601	HEM	CAC-C3C	2.85	1.55	1.47
2	A	601	HEM	FE-NB	2.50	2.02	1.94
2	A	601	HEM	FE-NC	2.47	2.03	1.95
2	A	601	HEM	FE-NA	2.36	2.02	1.95
4	A	603	SAM	C5-C6	2.30	1.47	1.41
4	A	603	SAM	C4-N3	2.29	1.38	1.34
4	A	603	SAM	CE-SD	-2.25	1.64	1.78
2	A	601	HEM	FE-ND	2.25	2.01	1.94
2	A	601	HEM	C2A-C3A	-2.19	1.33	1.38
2	A	601	HEM	CMB-C2B	2.14	1.55	1.50
4	A	603	SAM	C2-N3	2.14	1.37	1.33
2	B	601	HEM	FE-NB	2.14	2.01	1.94
4	A	603	SAM	C2'-C3'	-2.12	1.47	1.53
3	A	602	PLP	C2-N1	2.09	1.37	1.33
2	B	601	HEM	CMB-C2B	2.07	1.55	1.50
2	B	601	HEM	CMC-C2C	2.06	1.55	1.50
3	B	602	PLP	C2-N1	2.03	1.37	1.33
2	A	601	HEM	CMC-C2C	2.02	1.54	1.50
2	B	601	HEM	FE-NC	2.02	2.01	1.95

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	601	HEM	C2A-C3A	-2.00	1.33	1.38

All (45) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	603	SAM	C4-N9-C8	6.06	112.10	105.74
4	B	603	SAM	C5-C4-N3	-5.09	119.71	126.72
4	B	603	SAM	N3-C2-N1	-5.04	120.95	128.58
4	B	603	SAM	N3-C4-N9	5.02	135.70	127.17
4	A	603	SAM	N3-C2-N1	-4.53	121.72	128.58
3	B	602	PLP	C4A-C4-C5	4.17	125.24	120.94
4	B	603	SAM	C2-N3-C4	4.17	122.00	111.83
3	A	602	PLP	C4A-C4-C5	3.96	125.03	120.94
2	B	601	HEM	C4D-ND-C1D	3.91	109.84	105.21
4	B	603	SAM	N9-C8-N7	-3.85	108.47	113.94
2	A	601	HEM	C4D-ND-C1D	3.29	109.11	105.21
4	B	603	SAM	C4-C5-N7	-3.20	106.92	110.58
4	A	603	SAM	C4-N9-C8	3.15	109.04	105.74
4	A	603	SAM	O4'-C1'-N9	3.03	113.91	108.09
4	A	603	SAM	N9-C8-N7	-2.98	109.71	113.94
3	B	602	PLP	C6-C5-C4	2.79	120.38	118.10
4	B	603	SAM	C6-C5-N7	2.65	137.20	132.09
4	A	603	SAM	C2-N1-C6	2.62	123.04	118.73
4	A	603	SAM	N6-C6-N1	2.57	124.10	118.38
4	B	603	SAM	C5-N7-C8	2.53	107.42	103.45
2	A	601	HEM	CHD-C1D-ND	2.52	127.14	124.42
2	A	601	HEM	CAA-CBA-CGA	-2.51	107.00	113.67
3	A	602	PLP	C6-C5-C4	2.51	120.15	118.10
2	A	601	HEM	CAD-CBD-CGD	-2.47	107.11	113.67
2	B	601	HEM	CAD-CBD-CGD	-2.40	107.29	113.67
2	B	601	HEM	CHA-C4D-ND	2.35	127.27	124.37
4	B	603	SAM	OXT-C-O	-2.34	118.77	124.08
4	A	603	SAM	C2-N3-C4	2.33	117.51	111.83
3	A	602	PLP	C4A-C4-C3	-2.29	116.70	120.52
4	A	603	SAM	C5-C6-N6	-2.26	117.68	123.29
2	B	601	HEM	CHD-C1D-ND	2.25	126.84	124.42
4	A	603	SAM	O4'-C1'-C2'	-2.23	101.83	106.62
2	B	601	HEM	CAA-CBA-CGA	-2.23	107.75	113.67
4	A	603	SAM	C5-C4-N3	-2.15	123.75	126.72
2	A	601	HEM	C1B-NB-C4B	2.15	107.76	105.21
4	A	603	SAM	N3-C4-N9	2.14	130.81	127.17
4	A	603	SAM	C5-N7-C8	2.11	106.76	103.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	603	SAM	C2-N1-C6	2.10	122.19	118.73
2	B	601	HEM	CMD-C2D-C1D	2.07	128.26	125.03
2	A	601	HEM	C3B-C2B-C1B	2.06	107.96	106.41
3	B	602	PLP	C4A-C4-C3	-2.06	117.09	120.52
4	B	603	SAM	C4-N9-C1'	-2.05	121.83	126.63
2	A	601	HEM	CAD-C3D-C4D	2.04	128.26	124.70
3	A	602	PLP	C5-C6-N1	-2.04	120.52	123.83
2	A	601	HEM	CHA-C4D-ND	2.03	126.88	124.37

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	603	SAM	C4'-C5'-SD-CE
4	A	603	SAM	O4'-C4'-C5'-SD
4	B	603	SAM	O4'-C4'-C5'-SD
4	B	603	SAM	O-C-CA-N
4	A	603	SAM	OXT-C-CA-N
4	A	603	SAM	CB-CG-SD-C5'
4	B	603	SAM	CB-CG-SD-C5'
4	A	603	SAM	C-CA-CB-CG
4	A	603	SAM	CB-CG-SD-CE
4	B	603	SAM	OXT-C-CA-N
4	B	603	SAM	CB-CG-SD-CE

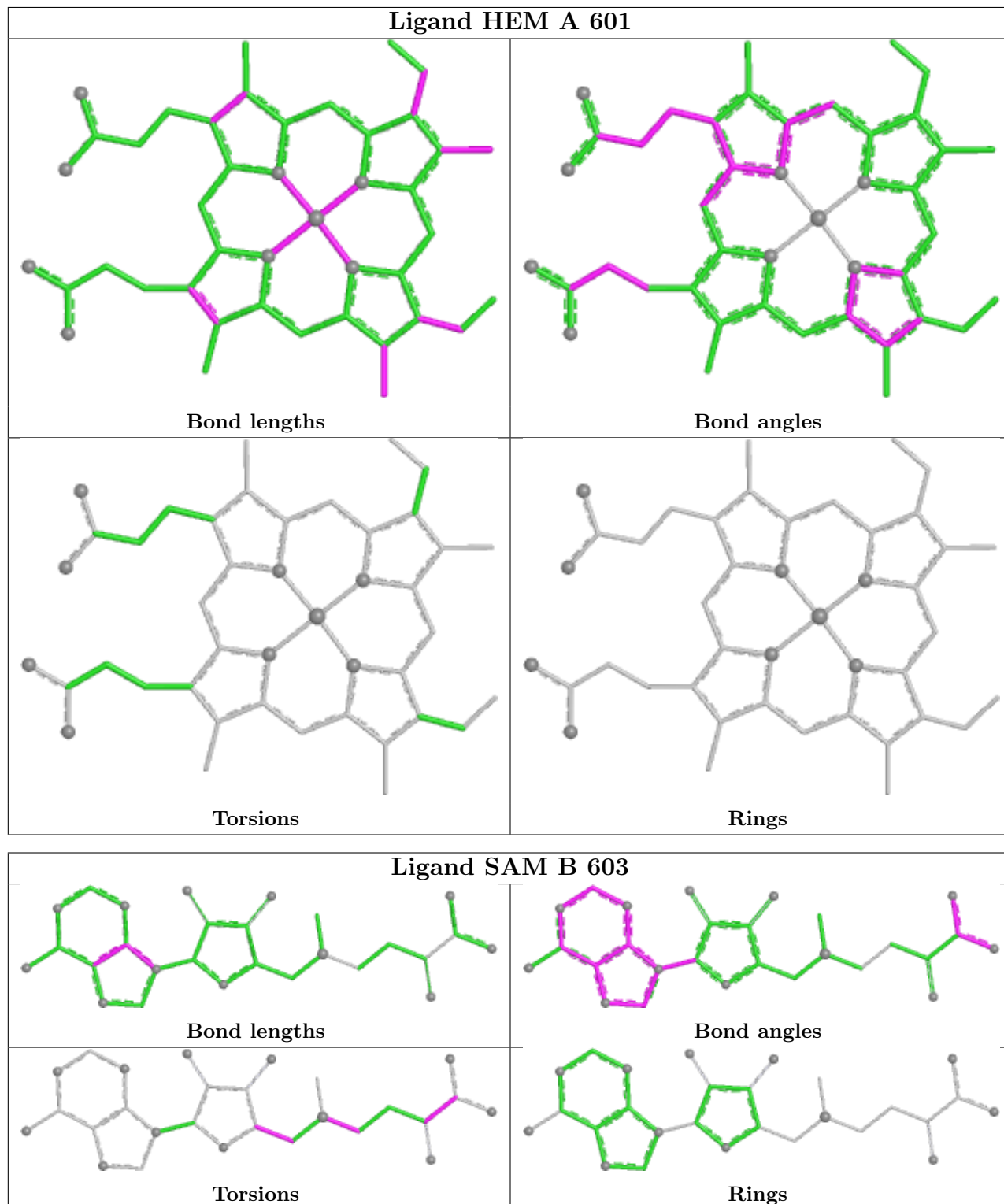
There are no ring outliers.

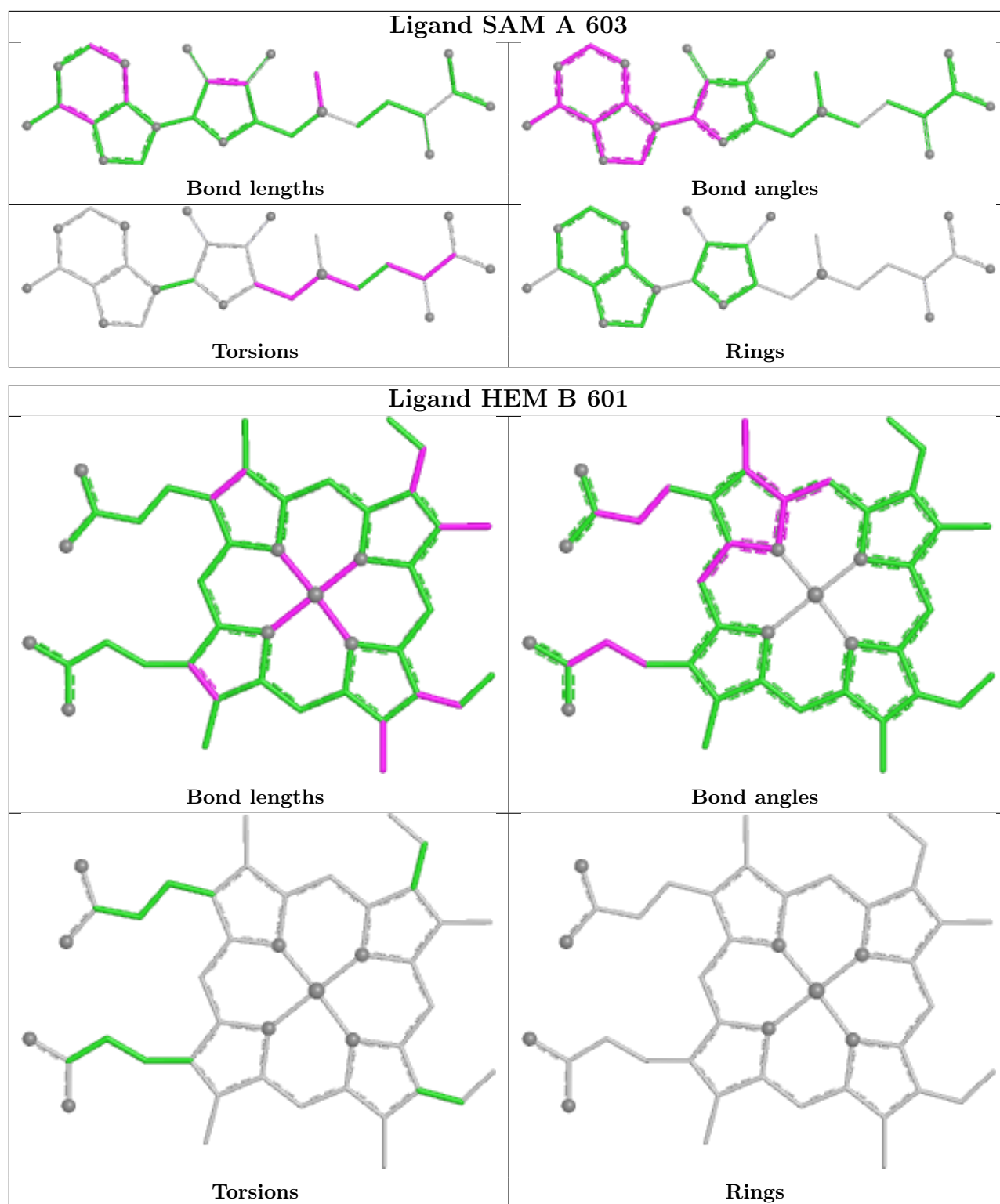
5 monomers are involved in 19 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	601	HEM	4	0
4	B	603	SAM	6	0
3	A	602	PLP	1	0
4	A	603	SAM	4	0
2	B	601	HEM	4	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

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	B	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	B	515:GLN	C	526:GLN	N	2.96

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	486/549 (88%)	0.16	2 (0%) 88 70	74, 95, 187, 212	0
1	B	489/549 (89%)	0.18	8 (1%) 70 42	74, 94, 222, 247	0
All	All	975/1098 (88%)	0.17	10 (1%) 79 53	74, 95, 211, 247	0

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	416	GLU	4.9
1	B	511	VAL	4.3
1	B	455	ILE	2.9
1	B	148	GLY	2.5
1	B	532	GLY	2.5
1	A	459	VAL	2.3
1	B	299	THR	2.1
1	B	120	ASP	2.1
1	A	388	ASP	2.0
1	B	297	GLU	2.0

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

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

6.3 Carbohydrates [i](#)

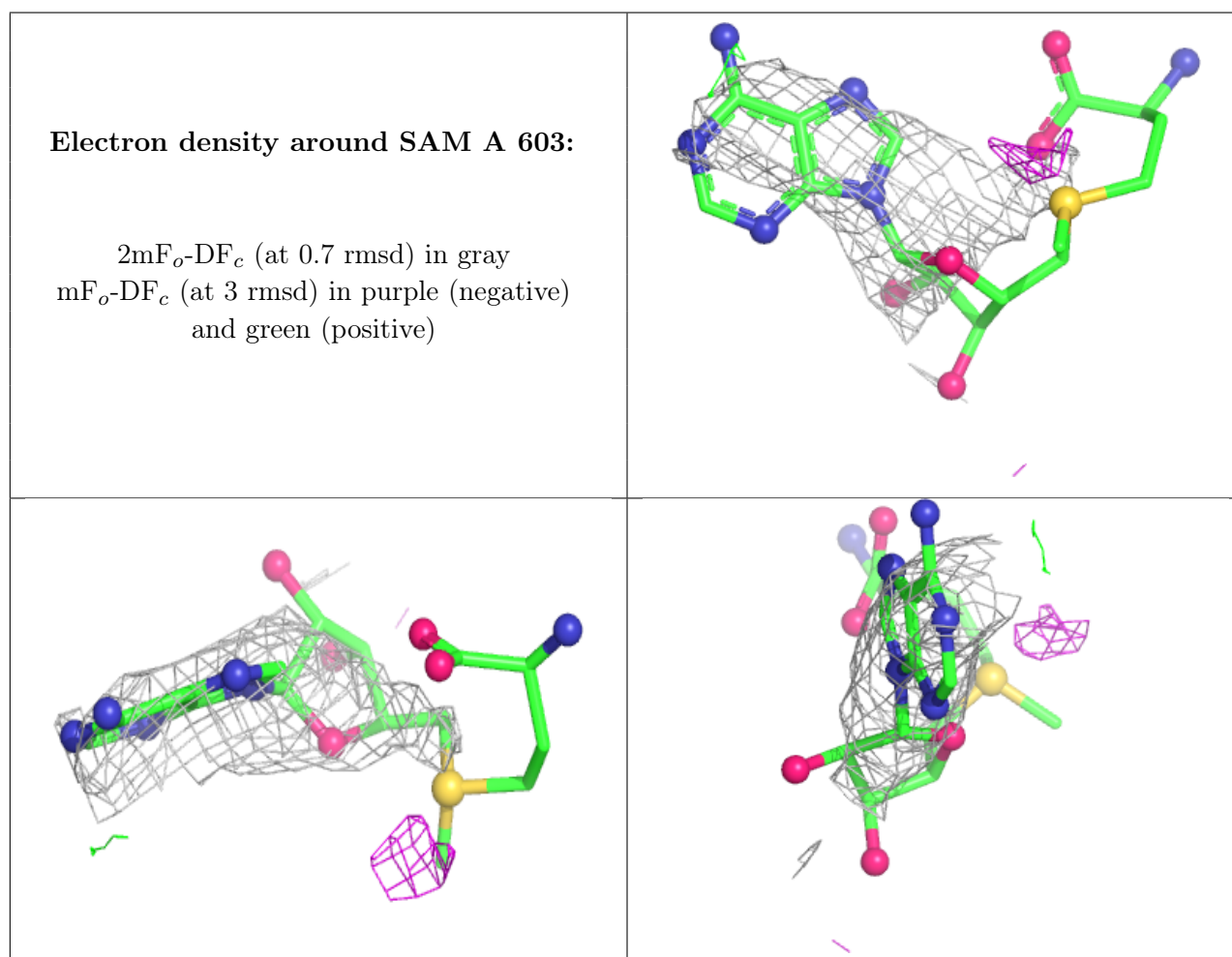
There are no oligosaccharides in this entry.

6.4 Ligands [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.

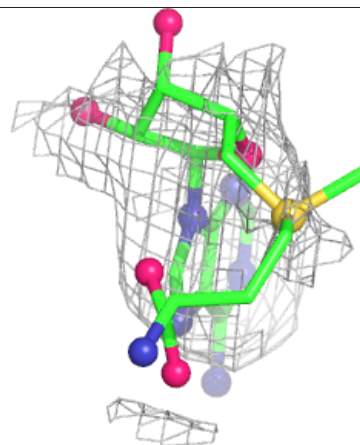
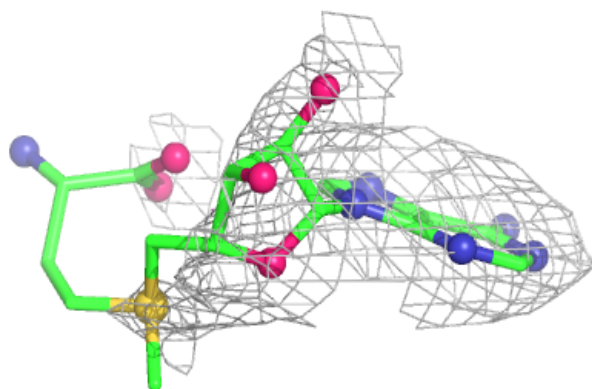
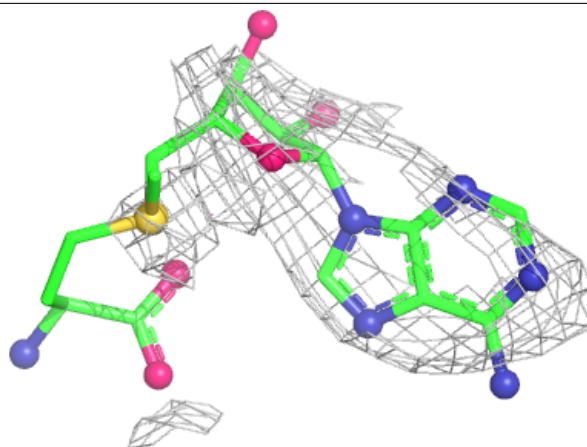
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	SAM	A	603	27/27	0.64	0.10	170,175,180,184	0
4	SAM	B	603	27/27	0.67	0.10	184,200,210,218	0
3	PLP	A	602	15/16	0.84	0.14	82,83,84,84	0
3	PLP	B	602	15/16	0.88	0.12	87,88,90,91	0
2	HEM	A	601	43/43	0.96	0.13	81,90,97,99	0
2	HEM	B	601	43/43	0.96	0.11	80,88,95,98	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.



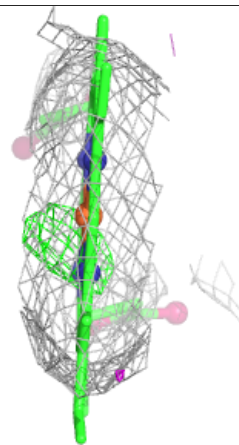
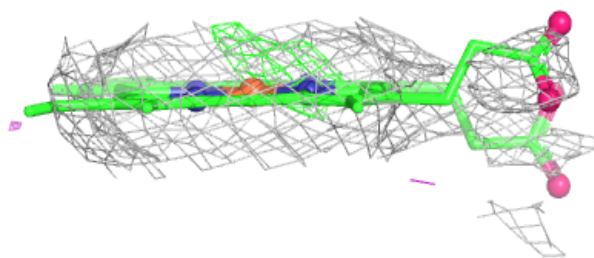
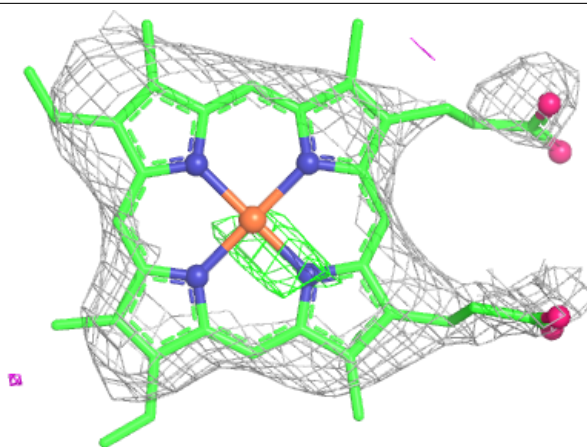
Electron density around SAM B 603:

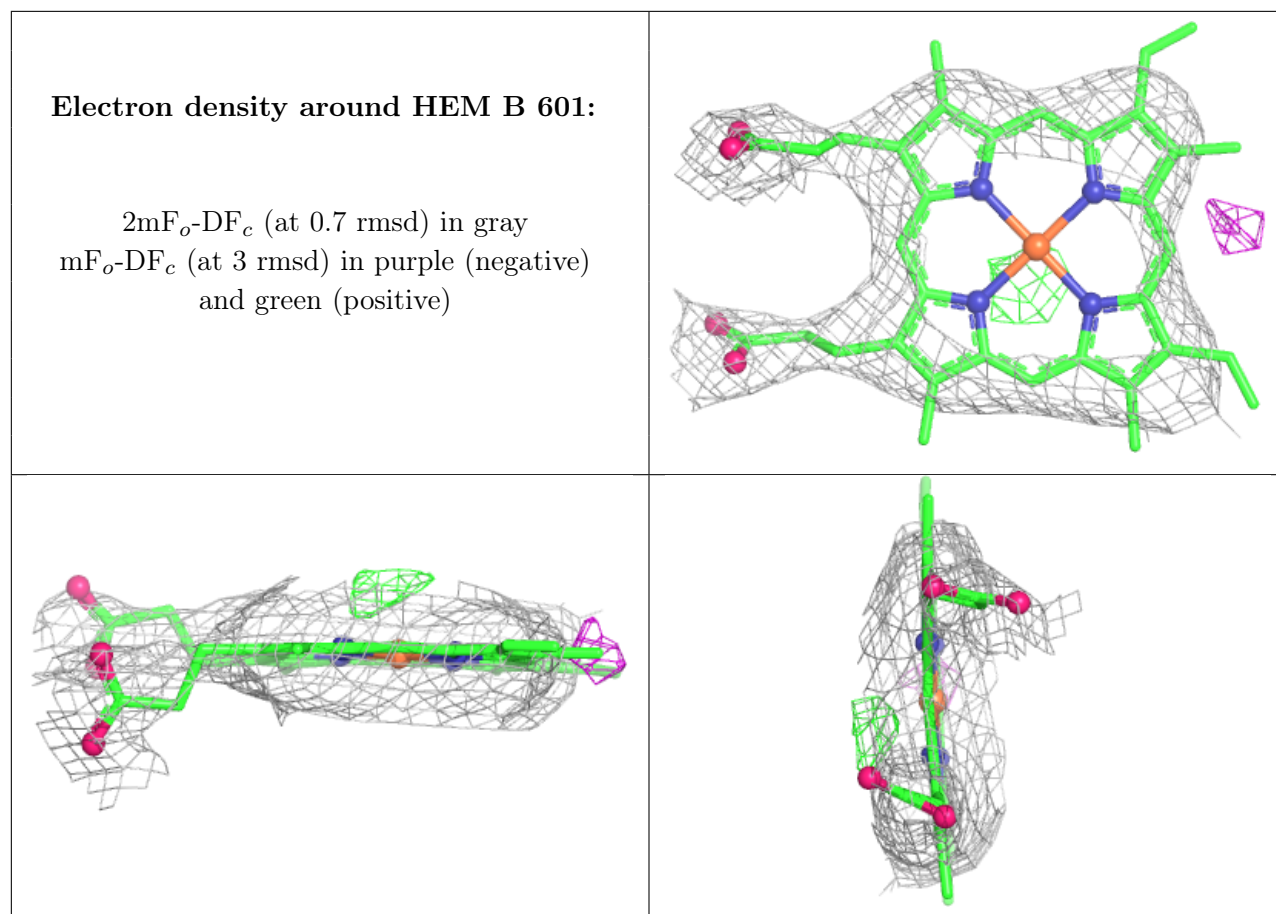
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around HEM A 601:

$2mF_o-DF_c$ (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.