



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 03:24 PM UTC

PDB ID : 5MVO / pdb_00005mvo
Title : FoxE P43212 crystal structure of Rhodopseudomonas ferrooxidans SW2 putative iron oxidase
Authors : Pereira, L.; Saraiva, I.H.; Oliveira, A.S.; Soares, C.; Gomes, R.O.; Frazao, C.
Deposited on : 2017-01-17
Resolution : 2.67 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

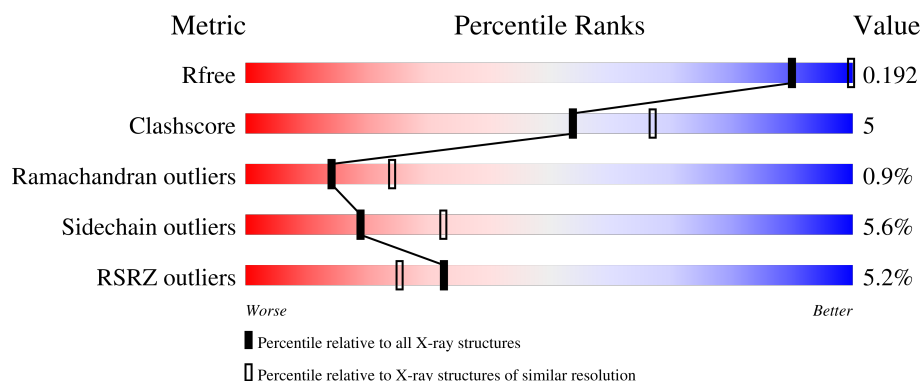
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

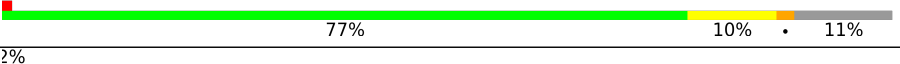
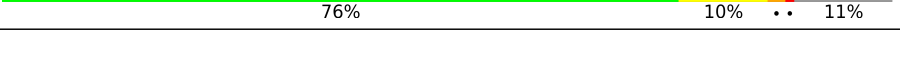
The reported resolution of this entry is 2.67 Å.

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	1110 (2.66-2.66)
Clashscore	190562	1141 (2.66-2.66)
Ramachandran outliers	187476	1126 (2.66-2.66)
Sidechain outliers	187428	1126 (2.66-2.66)
RSRZ outliers	180081	1110 (2.66-2.66)

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.

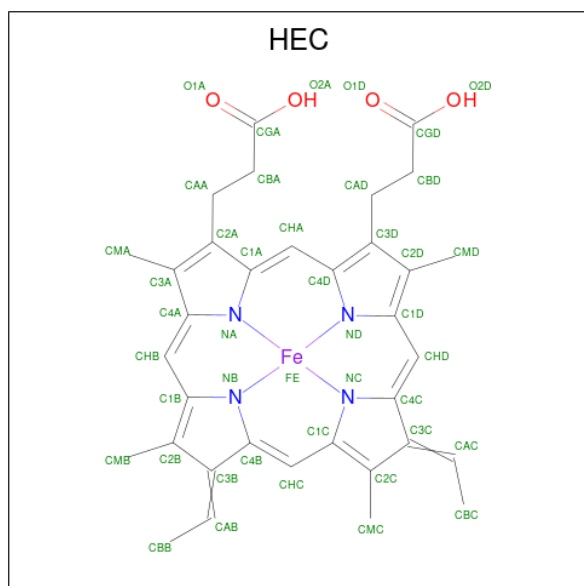
Mol	Chain	Length	Quality of chain
1	A	291	 11% 70% 14% .. 14%
1	B	291	 % 77% 10% • 11%
1	C	291	 2% 76% 10% .. 11%

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 FoxE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	251	Total 1927	C 1209	N 336	O 373	S 9	0	1	0
1	B	259	Total 2000	C 1251	N 347	O 393	S 9	0	3	0
1	C	259	Total 2001	C 1254	N 347	O 391	S 9	0	3	0

- Molecule 2 is HEME C (CCD ID: HEC) (formula: $\text{C}_{34}\text{H}_{34}\text{FeN}_4\text{O}_4$).



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

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

- Molecule 3 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	O P		
			5	4 1	0	0
3	A	1	Total	O P		
			5	4 1	0	0
3	B	1	Total	O P		
			5	4 1	0	0
3	B	1	Total	O P		
			5	4 1	0	0
3	C	1	Total	O P		
			5	4 1	0	0
3	C	1	Total	O P		
			5	4 1	0	0

- Molecule 4 is COPPER (II) ION (CCD ID: CU) (formula: Cu).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total 1	Cu 1	0	0
4	C	1	Total 1	Cu 1	0	0

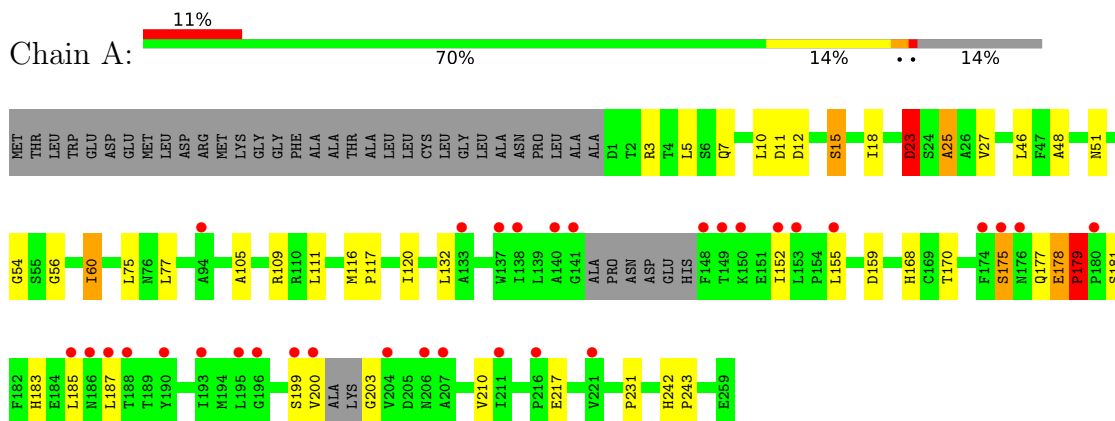
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	6	Total 6	O 6	0	0
5	B	20	Total 20	O 20	0	0
5	C	18	Total 18	O 18	0	0

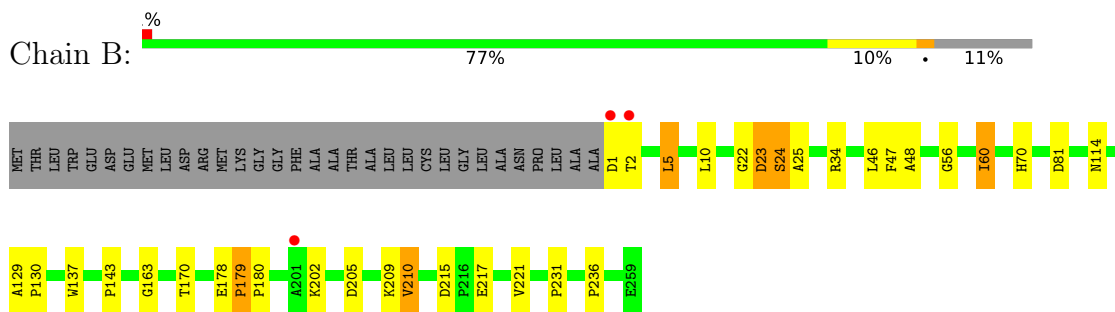
3 Residue-property plots [i](#)

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.

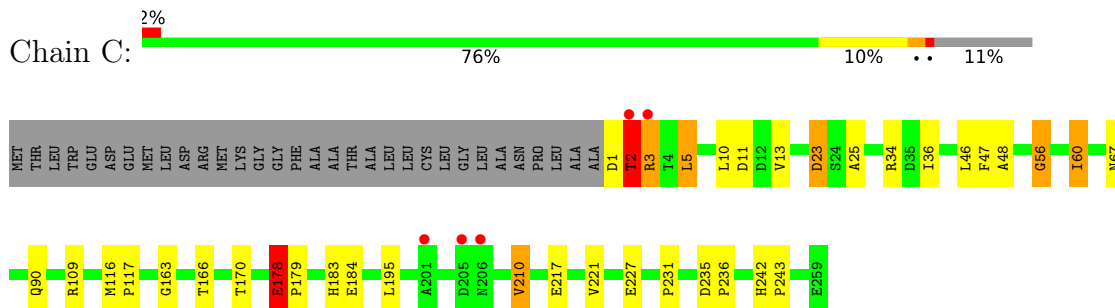
• Molecule 1: FoxE



• Molecule 1: FoxE



• Molecule 1: FoxE



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	116.85Å 116.85Å 191.13Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.85 – 2.67 49.85 – 2.67	Depositor EDS
% Data completeness (in resolution range)	99.3 (49.85-2.67) 99.2 (49.85-2.67)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.52 (at 2.65Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.174 , 0.174 0.192 , 0.192	Depositor DCC
R_{free} test set	2020 reflections (5.30%)	wwPDB-VP
Wilson B-factor (Å ²)	63.9	Xtriage
Anisotropy	0.131	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 59.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6262	wwPDB-VP
Average B, all atoms (Å ²)	76.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.70% 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: PO4, HEC, CU

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	0.57	0/1978	0.97	4/2693 (0.1%)
1	B	0.60	0/2055	1.05	9/2803 (0.3%)
1	C	0.61	0/2055	1.06	14/2803 (0.5%)
All	All	0.60	0/6088	1.03	27/8299 (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	1
1	C	0	2
All	All	0	4

There are no bond length outliers.

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	163	GLY	CA-C-N	8.73	128.75	119.76
1	C	163	GLY	C-N-CA	8.73	128.75	119.76
1	B	205	ASP	N-CA-C	-6.89	104.07	113.30
1	C	23	ASP	CA-C-N	6.83	134.00	121.70
1	C	23	ASP	C-N-CA	6.83	134.00	121.70
1	C	166	THR	CA-C-N	6.65	126.57	119.85
1	C	166	THR	C-N-CA	6.65	126.57	119.85
1	C	36	ILE	CA-C-N	6.23	126.44	119.90
1	C	36	ILE	C-N-CA	6.23	126.44	119.90
1	A	23	ASP	CA-C-N	6.18	132.83	121.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	23	ASP	C-N-CA	6.18	132.83	121.70
1	C	178	GLU	CA-C-N	5.81	126.36	120.38
1	C	178	GLU	C-N-CA	5.81	126.36	120.38
1	C	56	GLY	CA-C-N	5.62	125.94	119.92
1	C	56	GLY	C-N-CA	5.62	125.94	119.92
1	C	67	ASN	CA-C-N	5.54	125.25	119.82
1	C	67	ASN	C-N-CA	5.54	125.25	119.82
1	B	215	ASP	CA-C-N	5.48	124.82	118.85
1	B	215	ASP	C-N-CA	5.48	124.82	118.85
1	A	175	SER	N-CA-C	5.33	115.17	108.45
1	B	23[A]	ASP	CA-C-N	5.24	131.13	121.70
1	B	23[A]	ASP	C-N-CA	5.24	131.13	121.70
1	B	23[B]	ASP	CA-C-N	5.24	131.13	121.70
1	B	23[B]	ASP	C-N-CA	5.24	131.13	121.70
1	B	163	GLY	CA-C-N	5.20	125.21	119.90
1	B	163	GLY	C-N-CA	5.20	125.21	119.90
1	A	25	ALA	N-CA-C	5.19	116.94	111.28

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	178	GLU	Peptide
1	B	178	GLU	Peptide
1	C	178	GLU	Peptide
1	C	2[B]	THR	Peptide

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	1927	0	1870	25	0
1	B	2000	0	1928	14	1
1	C	2001	0	1935	17	0
2	A	86	0	61	1	0
2	B	86	0	60	3	0
2	C	86	0	60	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	10	0	0	0	0
3	B	10	0	0	0	0
3	C	10	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
5	A	6	0	0	0	0
5	B	20	0	0	0	0
5	C	18	0	0	0	0
All	All	6262	0	5914	60	1

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 5.

All (60) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:23:ASP:HA	1:C:25:ALA:H	1.18	1.08
1:A:23:ASP:HA	1:A:25:ALA:H	1.28	0.97
1:C:23:ASP:HA	1:C:25:ALA:N	1.99	0.75
1:A:23:ASP:HA	1:A:25:ALA:N	2.03	0.72
1:C:178:GLU:HB2	1:C:184:GLU:HG3	1.79	0.64
1:A:23:ASP:CA	1:A:25:ALA:H	2.07	0.63
1:B:210:VAL:HG13	1:B:221:VAL:HG23	1.86	0.57
1:A:3[B]:ARG:NE	1:A:11:ASP:OD2	2.31	0.57
1:C:48:ALA:HA	1:C:60:ILE:HG12	1.86	0.57
1:A:11:ASP:O	1:A:15:SER:OG	2.22	0.57
1:A:117:PRO:HD2	1:A:120:ILE:HD12	1.87	0.56
1:B:231:PRO:HD3	2:B:301:HEC:HBB2	1.88	0.56
1:A:3[B]:ARG:HG2	1:A:7:GLN:HB3	1.89	0.54
1:A:175:SER:OG	1:A:181:SER:HB2	2.08	0.54
1:A:48:ALA:HA	1:A:60:ILE:HG12	1.89	0.54
1:C:1[A]:ASP:O	1:C:2[A]:THR:HG23	2.08	0.53
1:C:210:VAL:HG13	1:C:221:VAL:HG23	1.90	0.53
1:C:56:GLY:HA3	1:C:231:PRO:HG2	1.92	0.52
1:A:12:ASP:HB3	1:A:18:ILE:HG13	1.95	0.49
1:A:152:ILE:O	1:A:155:LEU:HD13	2.13	0.49
1:A:159:ASP:OD1	1:A:168:HIS:ND1	2.32	0.49
1:A:185:LEU:HD21	1:A:187:LEU:HD21	1.94	0.49
1:B:48:ALA:HA	1:B:60:ILE:HG12	1.95	0.48
1:A:178:GLU:HG2	1:A:179:PRO:HD2	1.95	0.48
1:A:105:ALA:O	1:A:109:ARG:HG3	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:301:HEC:HMC3	2:C:301:HEC:HBC3	1.96	0.48
1:A:75:LEU:HD21	1:A:77:LEU:HD21	1.96	0.47
1:C:109:ARG:NH1	1:C:227:GLU:OE2	2.49	0.46
1:A:132:LEU:HD23	1:A:132:LEU:HA	1.81	0.45
1:C:183:HIS:ND1	2:C:302:HEC:O2A	2.49	0.45
1:A:177:GLN:HG2	1:A:200:VAL:HG21	1.99	0.45
2:B:301:HEC:HBC3	2:B:301:HEC:HMC3	1.99	0.45
1:A:199:SER:C	1:A:203:GLY:HA3	2.42	0.44
1:B:129:ALA:HB3	1:B:130:PRO:HD3	1.99	0.44
1:B:56:GLY:HA3	1:B:231:PRO:HG2	1.99	0.44
1:C:47:PHE:O	1:C:60:ILE:HG23	2.17	0.44
1:A:242:HIS:ND1	1:A:243:PRO:HD2	2.34	0.43
1:C:235:ASP:HA	1:C:236:PRO:HD3	1.90	0.43
1:C:5:LEU:HD12	1:C:5:LEU:HA	1.77	0.43
1:C:1[B]:ASP:O	1:C:2[B]:THR:OG1	2.33	0.43
1:A:178:GLU:HG2	1:A:179:PRO:CD	2.48	0.43
1:A:51:ASN:HA	1:A:54:GLY:O	2.19	0.43
1:A:56:GLY:HA3	1:A:231:PRO:HG2	2.01	0.43
1:B:114:ASN:OD1	1:B:236:PRO:HA	2.20	0.42
1:B:5:LEU:HD12	1:B:5:LEU:HA	1.70	0.42
1:B:137:TRP:CD1	1:B:143:PRO:HD2	2.54	0.42
1:C:116:MET:HA	1:C:117:PRO:C	2.45	0.42
1:A:111:LEU:HD23	1:A:111:LEU:HA	1.93	0.42
1:B:22:GLY:O	1:B:24:SER:HB3	2.19	0.42
1:C:1[A]:ASP:HB3	1:C:3:ARG:HG2	2.01	0.41
2:B:302:HEC:HMC1	2:B:302:HEC:HBC3	2.02	0.41
1:C:242:HIS:ND1	1:C:243:PRO:HD2	2.35	0.41
1:C:3:ARG:NH2	1:C:11:ASP:OD2	2.53	0.41
1:B:179:PRO:HA	1:B:180:PRO:HA	1.65	0.41
2:A:301:HEC:HBC3	2:A:301:HEC:HMC3	2.03	0.41
1:B:47:PHE:O	1:B:60:ILE:HG23	2.20	0.41
1:A:116:MET:HA	1:A:117:PRO:C	2.46	0.40

All (1) 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:1[B]:ASP:OD1	1:B:70:HIS:NE2[3_544]	2.04	0.16

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	246/291 (84%)	240 (98%)	5 (2%)	1 (0%)	30	45
1	B	259/291 (89%)	246 (95%)	9 (4%)	4 (2%)	8	13
1	C	259/291 (89%)	249 (96%)	6 (2%)	4 (2%)	8	13
All	All	764/873 (88%)	735 (96%)	20 (3%)	9 (1%)	14	17

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	179	PRO
1	B	179	PRO
1	B	202	LYS
1	C	2[A]	THR
1	C	2[B]	THR
1	C	3	ARG
1	C	179	PRO
1	B	2[A]	THR
1	B	2[B]	THR

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	211/239 (88%)	199 (94%)	12 (6%)	18	32
1	B	219/239 (92%)	206 (94%)	13 (6%)	18	30
1	C	197/239 (82%)	185 (94%)	12 (6%)	17	30

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
All	All	627/717 (87%)	590 (94%)	37 (6%)	19	30

All (37) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	LEU
1	A	10	LEU
1	A	15	SER
1	A	23	ASP
1	A	27	VAL
1	A	46	LEU
1	A	60	ILE
1	A	170	THR
1	A	179	PRO
1	A	183	HIS
1	A	210	VAL
1	A	217	GLU
1	B	5	LEU
1	B	10	LEU
1	B	23[A]	ASP
1	B	23[B]	ASP
1	B	24	SER
1	B	34	ARG
1	B	46	LEU
1	B	60	ILE
1	B	81	ASP
1	B	170	THR
1	B	209	LYS
1	B	210	VAL
1	B	217	GLU
1	C	5	LEU
1	C	10[A]	LEU
1	C	10[B]	LEU
1	C	13	VAL
1	C	34	ARG
1	C	46	LEU
1	C	60	ILE
1	C	90	GLN
1	C	170	THR
1	C	195	LEU
1	C	210	VAL
1	C	217	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	90	GLN
1	A	246	GLN
1	B	8	GLN
1	B	104	ASN
1	B	147	HIS
1	C	246	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 ⓘ

Of 14 ligands modelled in this entry, 2 are monoatomic - leaving 12 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
2	HEC	B	302	1	46,50,50	2.05	7 (15%)	58,82,82	1.07	4 (6%)
3	PO4	A	303	-	4,4,4	0.94	0	6,6,6	0.63	0
2	HEC	C	301	1	46,50,50	2.08	8 (17%)	58,82,82	1.12	3 (5%)
2	HEC	A	301	1	46,50,50	2.00	7 (15%)	58,82,82	1.18	5 (8%)
3	PO4	A	304	-	4,4,4	0.94	0	6,6,6	0.58	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	PO4	C	303	-	4,4,4	0.84	0	6,6,6	0.53	0
3	PO4	B	303	-	4,4,4	0.96	0	6,6,6	0.76	0
2	HEC	C	302	1	46,50,50	2.05	10 (21%)	58,82,82	1.19	4 (6%)
3	PO4	C	304	-	4,4,4	0.92	0	6,6,6	0.71	0
3	PO4	B	304	-	4,4,4	0.87	0	6,6,6	0.90	0
2	HEC	A	302	1	46,50,50	2.09	9 (19%)	58,82,82	1.12	3 (5%)
2	HEC	B	301	1	46,50,50	1.99	7 (15%)	58,82,82	1.25	5 (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	HEC	B	302	1	-	4/14/54/54	-
2	HEC	C	301	1	-	7/14/54/54	-
2	HEC	A	301	1	-	6/14/54/54	-
2	HEC	C	302	1	-	4/14/54/54	-
2	HEC	A	302	1	-	2/14/54/54	-
2	HEC	B	301	1	-	6/14/54/54	-

All (48) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	302	HEC	CAB-C3B	6.54	1.56	1.35
2	A	302	HEC	CAC-C3C	6.46	1.56	1.35
2	B	301	HEC	CAB-C3B	6.39	1.55	1.35
2	B	302	HEC	CAC-C3C	6.38	1.55	1.35
2	C	302	HEC	CAB-C3B	6.34	1.55	1.35
2	C	301	HEC	CAC-C3C	6.34	1.55	1.35
2	C	301	HEC	CAB-C3B	6.33	1.55	1.35
2	A	301	HEC	CAC-C3C	6.30	1.55	1.35
2	C	302	HEC	CAC-C3C	6.26	1.55	1.35
2	B	301	HEC	CAC-C3C	6.14	1.55	1.35
2	B	302	HEC	CAB-C3B	5.99	1.54	1.35
2	C	301	HEC	C3D-C2D	5.88	1.54	1.38
2	A	301	HEC	CAB-C3B	5.73	1.53	1.35
2	A	302	HEC	C3D-C2D	5.62	1.53	1.38
2	C	302	HEC	C3D-C2D	5.58	1.53	1.38
2	B	302	HEC	C3D-C2D	5.44	1.53	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	301	HEC	C3D-C2D	5.27	1.52	1.38
2	A	301	HEC	C3D-C2D	5.01	1.52	1.38
2	A	301	HEC	CBB-CAB	-4.31	1.33	1.49
2	A	301	HEC	CBC-CAC	-4.24	1.33	1.49
2	C	302	HEC	CBC-CAC	-4.23	1.33	1.49
2	A	302	HEC	CBB-CAB	-4.15	1.34	1.49
2	B	302	HEC	CBB-CAB	-4.13	1.34	1.49
2	B	301	HEC	CBC-CAC	-4.12	1.34	1.49
2	A	302	HEC	CBC-CAC	-4.09	1.34	1.49
2	C	301	HEC	CBB-CAB	-4.09	1.34	1.49
2	B	302	HEC	CBC-CAC	-4.09	1.34	1.49
2	C	301	HEC	CBC-CAC	-4.05	1.34	1.49
2	B	301	HEC	CBB-CAB	-3.97	1.34	1.49
2	C	302	HEC	CBB-CAB	-3.91	1.35	1.49
2	A	301	HEC	CMC-C2C	2.53	1.56	1.50
2	A	302	HEC	CAA-C2A	2.36	1.57	1.51
2	B	302	HEC	C3C-C2C	-2.35	1.33	1.41
2	C	302	HEC	C3B-C2B	-2.29	1.33	1.41
2	C	301	HEC	CMD-C2D	2.28	1.55	1.50
2	A	301	HEC	C3B-C2B	-2.28	1.33	1.41
2	A	302	HEC	CMB-C2B	2.26	1.55	1.50
2	B	301	HEC	CMC-C2C	2.25	1.55	1.50
2	C	301	HEC	C3C-C2C	-2.24	1.33	1.41
2	B	302	HEC	CMB-C2B	2.22	1.55	1.50
2	C	302	HEC	CMB-C2B	2.19	1.55	1.50
2	C	302	HEC	CMC-C2C	2.16	1.55	1.50
2	C	301	HEC	C3B-C2B	-2.14	1.33	1.41
2	C	302	HEC	CMD-C2D	2.14	1.55	1.50
2	A	302	HEC	CMC-C2C	2.11	1.55	1.50
2	C	302	HEC	C3C-C2C	-2.08	1.34	1.41
2	B	301	HEC	C3B-C2B	-2.03	1.34	1.41
2	A	302	HEC	C3C-C2C	-2.01	1.34	1.41

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	301	HEC	CBB-CAB-C3B	-4.42	118.60	127.43
2	B	301	HEC	CBB-CAB-C3B	-4.27	118.89	127.43
2	C	301	HEC	CBB-CAB-C3B	-3.64	120.16	127.43
2	B	302	HEC	CBB-CAB-C3B	-3.63	120.18	127.43
2	C	302	HEC	CBC-CAC-C3C	-3.62	120.20	127.43
2	C	302	HEC	CBB-CAB-C3B	-3.49	120.46	127.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	302	HEC	C4D-ND-C1D	3.39	111.34	105.82
2	C	302	HEC	C4D-ND-C1D	3.21	111.05	105.82
2	C	301	HEC	C4D-ND-C1D	3.20	111.04	105.82
2	A	301	HEC	C4D-ND-C1D	3.06	110.81	105.82
2	B	301	HEC	C4D-ND-C1D	3.06	110.80	105.82
2	A	301	HEC	CAD-C3D-C4D	2.55	129.93	124.94
2	B	302	HEC	C4D-ND-C1D	2.52	109.92	105.82
2	C	302	HEC	CHB-C4A-NA	2.51	127.18	124.45
2	B	301	HEC	CAD-C3D-C4D	2.43	129.68	124.94
2	B	302	HEC	CAD-C3D-C4D	2.24	129.32	124.94
2	B	302	HEC	CHC-C4B-NB	2.13	126.78	124.45
2	B	301	HEC	O2D-CGD-CBD	2.13	120.73	114.00
2	C	301	HEC	CAD-C3D-C4D	2.11	129.06	124.94
2	A	302	HEC	C4A-C3A-C2A	2.10	110.09	106.97
2	A	301	HEC	CBC-CAC-C3C	-2.09	123.26	127.43
2	A	301	HEC	CMD-C2D-C1D	2.07	128.57	125.42
2	A	302	HEC	CBB-CAB-C3B	-2.02	123.40	127.43
2	B	301	HEC	CBC-CAC-C3C	-2.02	123.41	127.43

There are no chirality outliers.

All (29) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	301	HEC	C2B-C3B-CAB-CBB
2	A	301	HEC	C4B-C3B-CAB-CBB
2	A	301	HEC	C2C-C3C-CAC-CBC
2	A	301	HEC	C4C-C3C-CAC-CBC
2	A	302	HEC	C2B-C3B-CAB-CBB
2	A	302	HEC	C4B-C3B-CAB-CBB
2	B	301	HEC	C2B-C3B-CAB-CBB
2	B	301	HEC	C4B-C3B-CAB-CBB
2	B	301	HEC	C2C-C3C-CAC-CBC
2	B	301	HEC	C4C-C3C-CAC-CBC
2	B	302	HEC	C2B-C3B-CAB-CBB
2	B	302	HEC	C4B-C3B-CAB-CBB
2	B	302	HEC	C2C-C3C-CAC-CBC
2	B	302	HEC	C4C-C3C-CAC-CBC
2	C	301	HEC	C2B-C3B-CAB-CBB
2	C	301	HEC	C4B-C3B-CAB-CBB
2	C	301	HEC	C2C-C3C-CAC-CBC
2	C	301	HEC	C4C-C3C-CAC-CBC
2	C	302	HEC	C2B-C3B-CAB-CBB

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Mol	Chain	Res	Type	Atoms
2	C	302	HEC	C4B-C3B-CAB-CBB
2	C	302	HEC	C2C-C3C-CAC-CBC
2	C	302	HEC	C4C-C3C-CAC-CBC
2	C	301	HEC	C1A-C2A-CAA-CBA
2	B	301	HEC	CAD-CBD-CGD-O2D
2	B	301	HEC	CAD-CBD-CGD-O1D
2	A	301	HEC	CAD-CBD-CGD-O2D
2	A	301	HEC	CAD-CBD-CGD-O1D
2	C	301	HEC	C3A-C2A-CAA-CBA
2	C	301	HEC	CAD-CBD-CGD-O2D

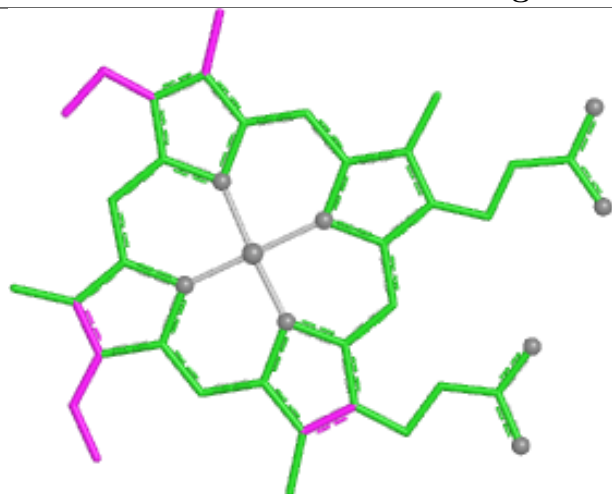
There are no ring outliers.

5 monomers are involved in 6 short contacts:

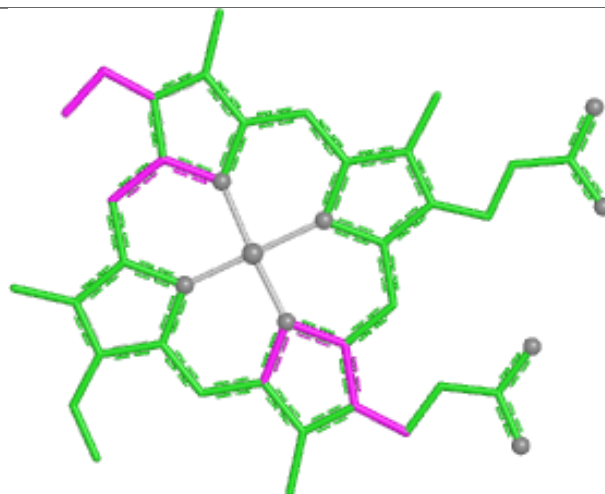
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	302	HEC	1	0
2	C	301	HEC	1	0
2	A	301	HEC	1	0
2	C	302	HEC	1	0
2	B	301	HEC	2	0

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.

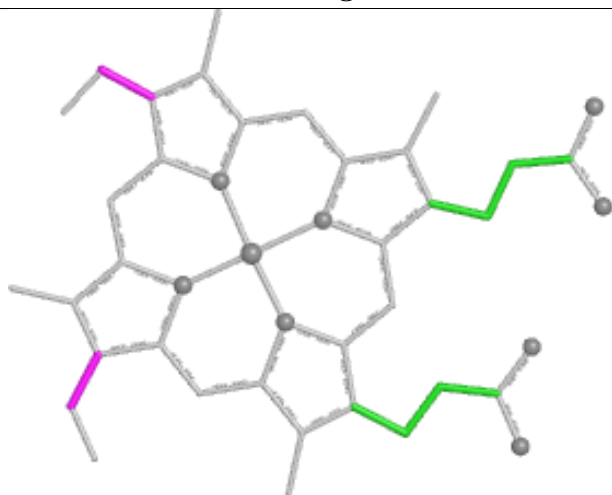
Ligand HEC B 302



Bond lengths



Bond angles

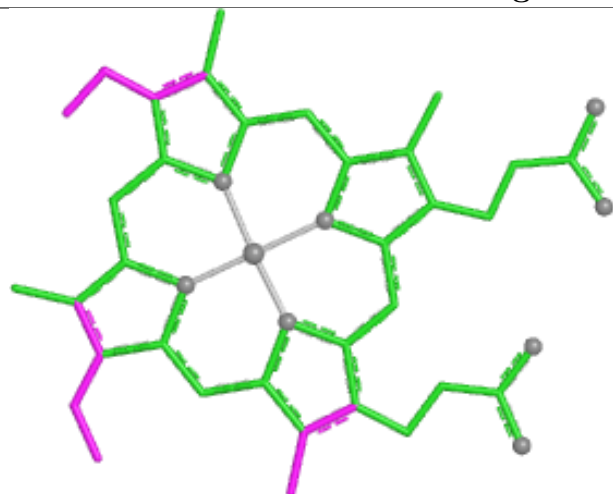


Torsions

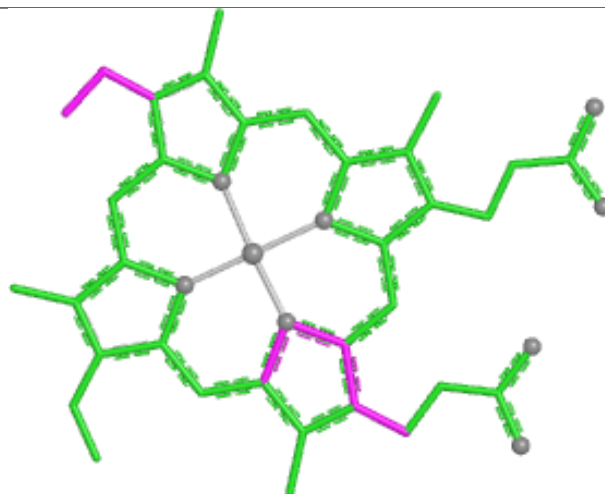


Rings

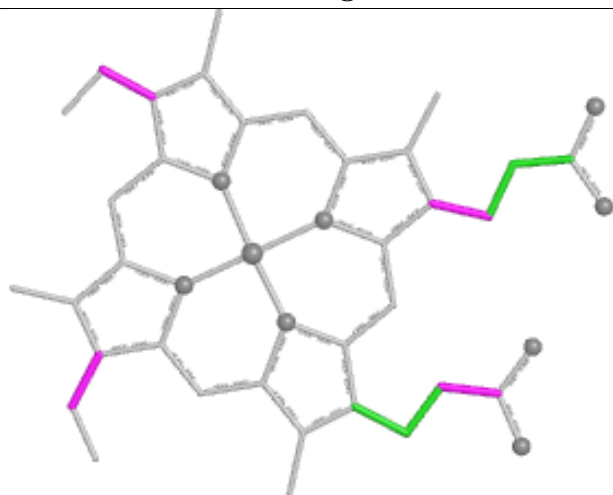
Ligand HEC C 301



Bond lengths



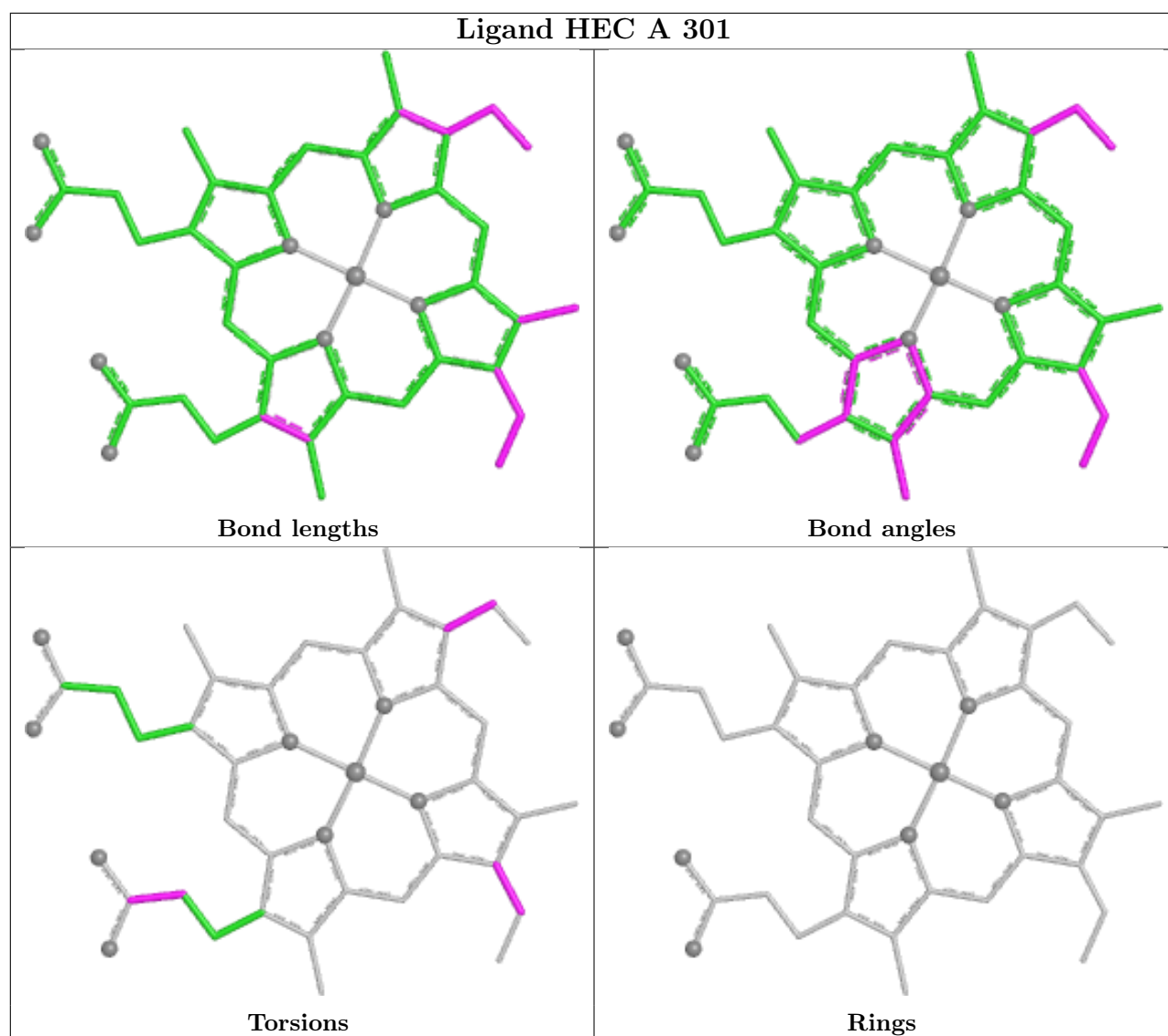
Bond angles



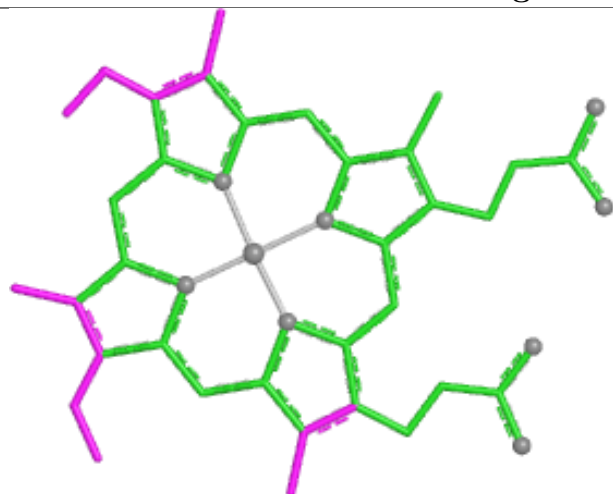
Torsions



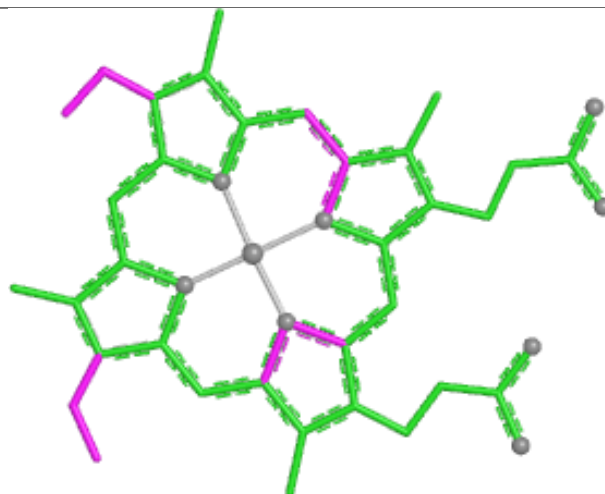
Rings



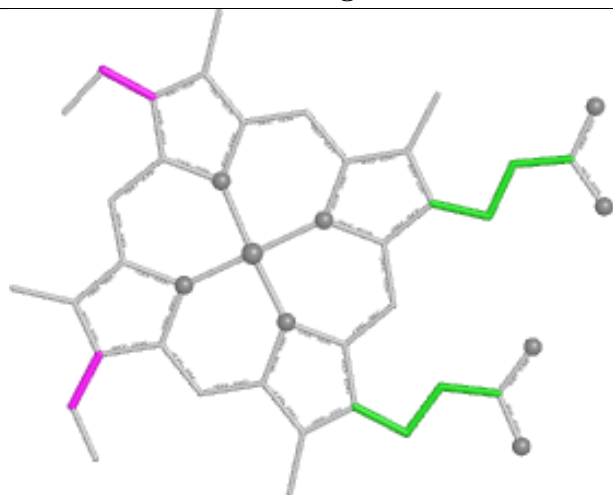
Ligand HEC C 302



Bond lengths



Bond angles

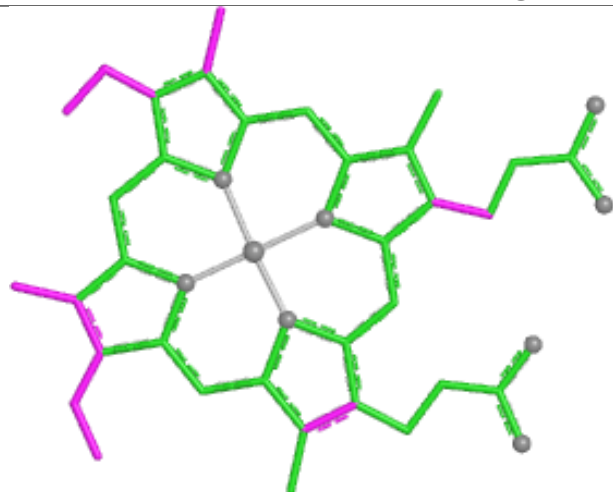


Torsions

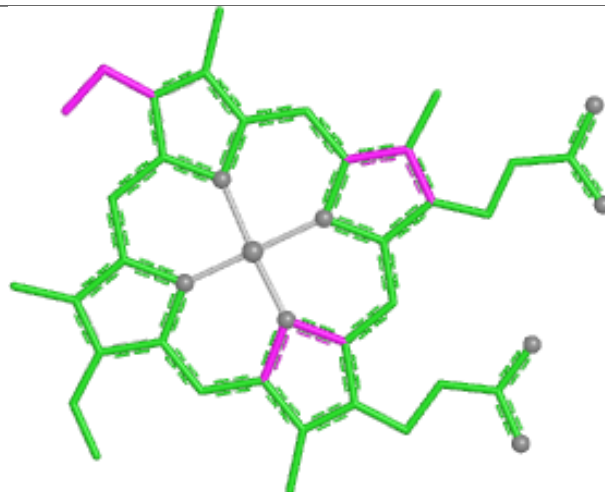


Rings

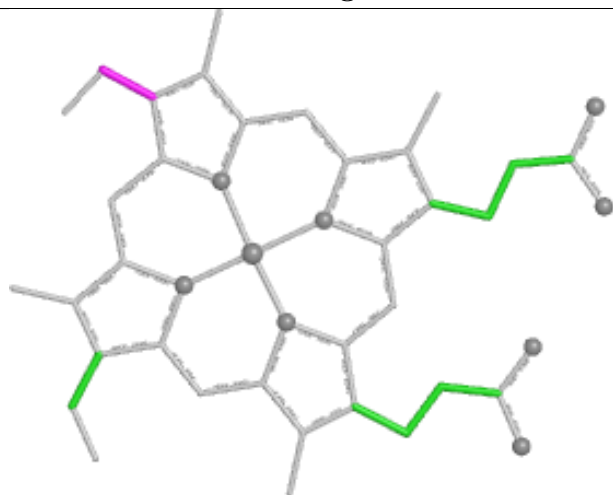
Ligand HEC A 302



Bond lengths



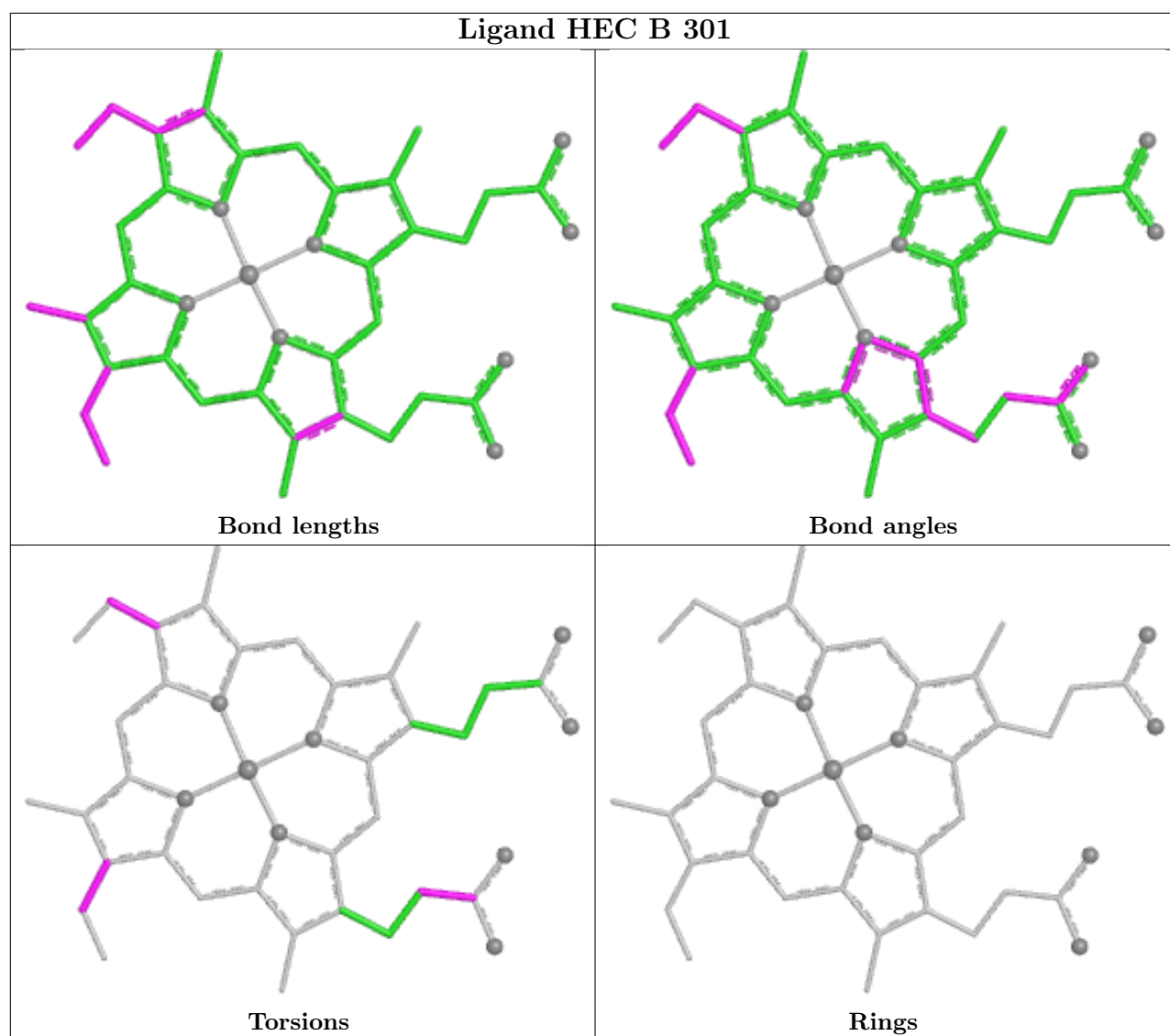
Bond angles



Torsions



Rings



5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	251/291 (86%)	0.53	32 (12%) 8 6	43, 88, 170, 197	1 (0%)
1	B	259/291 (89%)	-0.32	3 (1%) 76 72	40, 61, 99, 133	3 (1%)
1	C	259/291 (89%)	-0.33	5 (1%) 66 61	41, 61, 96, 132	3 (1%)
All	All	769/873 (88%)	-0.04	40 (5%) 33 25	40, 66, 157, 197	7 (0%)

All (40) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	148	PHE	5.2
1	C	206	ASN	5.0
1	A	196	GLY	4.4
1	A	149	THR	4.0
1	A	137	TRP	3.7
1	C	2[A]	THR	3.5
1	A	190	TYR	3.3
1	A	140	ALA	3.3
1	A	204	VAL	3.3
1	B	201	ALA	3.2
1	A	185	LEU	3.1
1	A	152	ILE	3.0
1	A	187	LEU	3.0
1	A	195	LEU	2.9
1	A	150	LYS	2.9
1	C	205	ASP	2.8
1	B	1[A]	ASP	2.8
1	A	141	GLY	2.7
1	A	180	PRO	2.7
1	A	199	SER	2.6
1	A	216	PRO	2.6
1	B	2[A]	THR	2.6
1	A	200	VAL	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	138	ILE	2.5
1	A	153	LEU	2.5
1	A	174	PHE	2.5
1	A	176	ASN	2.4
1	A	133	ALA	2.4
1	A	193	ILE	2.4
1	A	221	VAL	2.4
1	A	186	ASN	2.3
1	A	188	THR	2.3
1	C	3	ARG	2.3
1	A	207	ALA	2.3
1	A	206	ASN	2.3
1	C	201	ALA	2.2
1	A	211	ILE	2.2
1	A	175	SER	2.2
1	A	94	ALA	2.1
1	A	155	LEU	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.

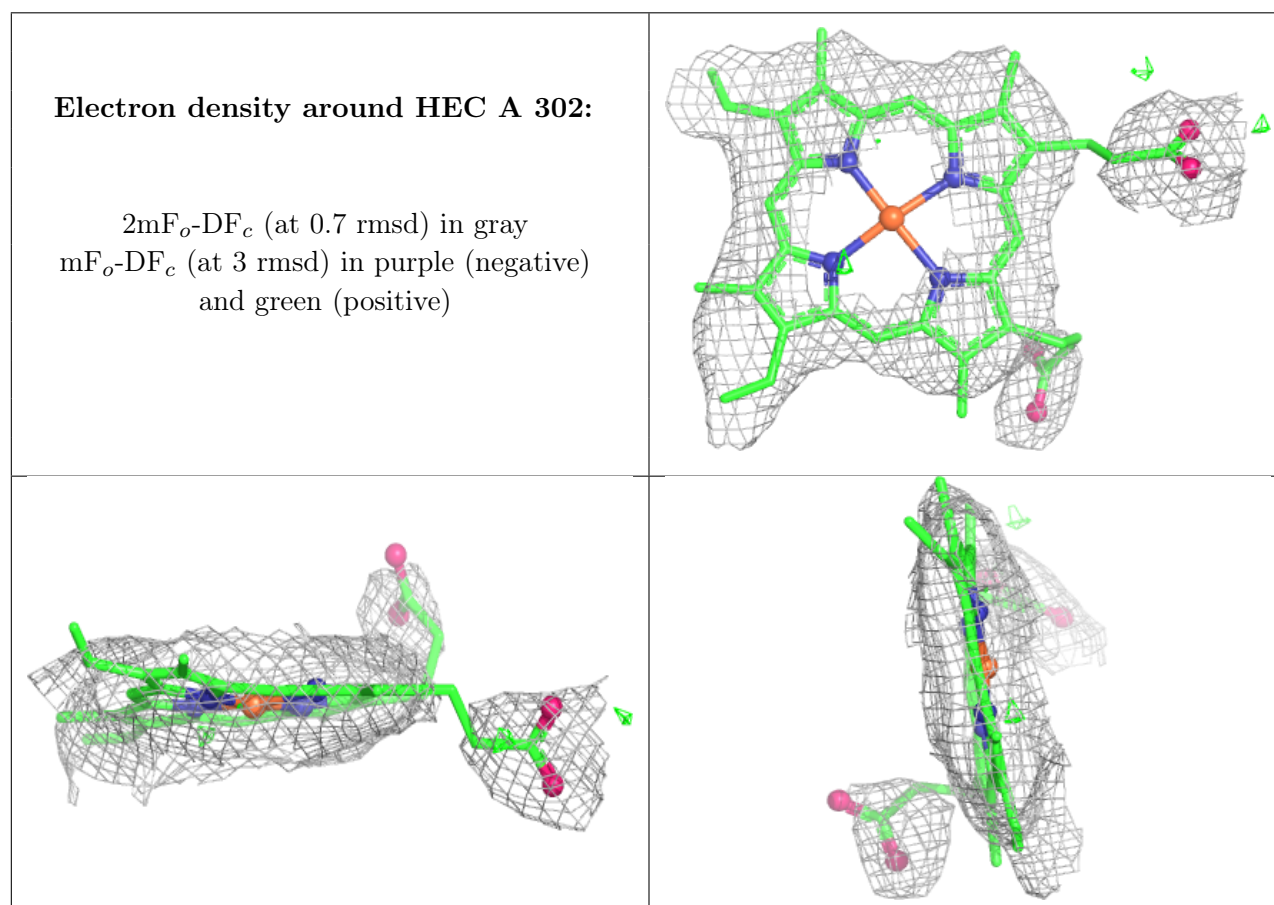
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	PO4	A	304	5/5	0.77	0.12	125,126,128,135	0
3	PO4	A	303	5/5	0.88	0.08	104,106,116,120	0
3	PO4	B	304	5/5	0.89	0.10	85,96,100,108	0
3	PO4	C	303	5/5	0.90	0.07	91,95,99,106	0
3	PO4	C	304	5/5	0.92	0.10	73,89,94,96	0
4	CU	C	305	1/1	0.96	0.07	122,122,122,122	0

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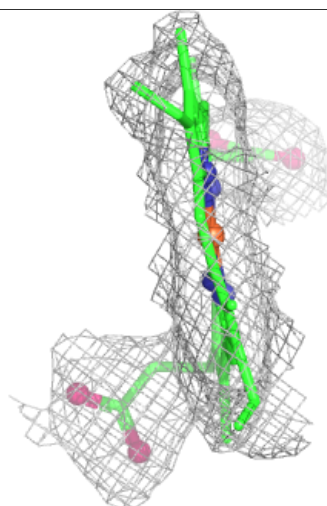
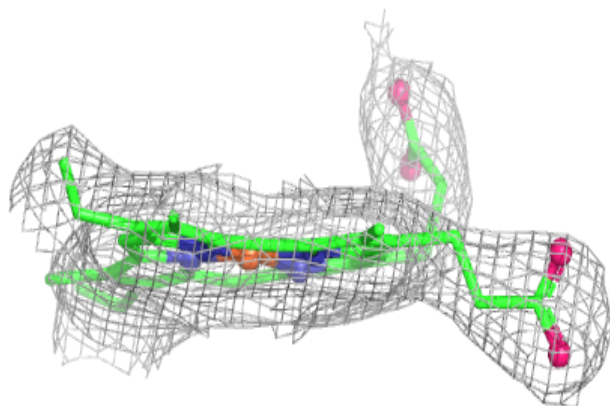
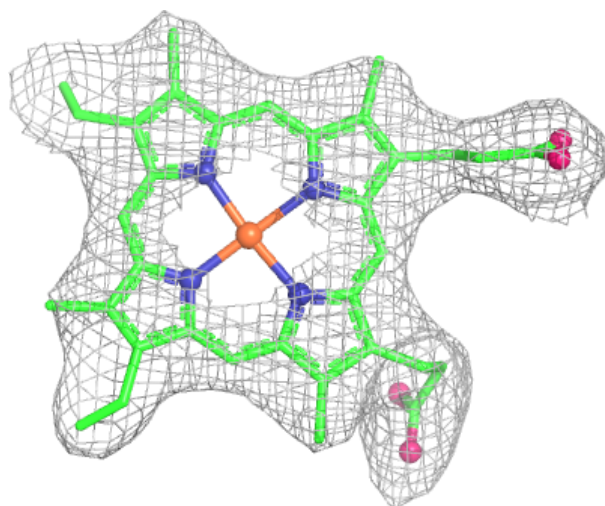
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	HEC	A	302	43/43	0.97	0.10	103,115,136,164	0
3	PO4	B	303	5/5	0.97	0.06	59,70,72,82	0
2	HEC	B	302	43/43	0.99	0.06	42,52,61,69	0
2	HEC	C	301	43/43	0.99	0.07	34,46,62,85	0
2	HEC	C	302	43/43	0.99	0.06	37,49,64,74	0
2	HEC	A	301	43/43	0.99	0.07	44,62,74,76	0
4	CU	B	305	1/1	0.99	0.05	90,90,90,90	0
2	HEC	B	301	43/43	0.99	0.06	39,52,60,69	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.



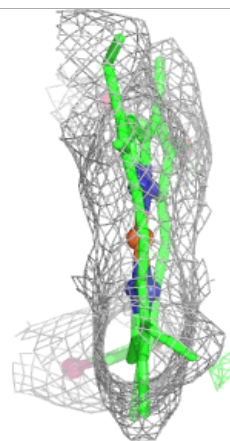
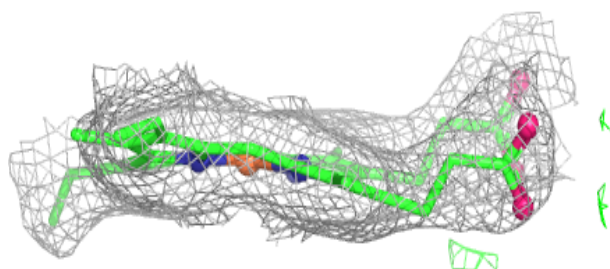
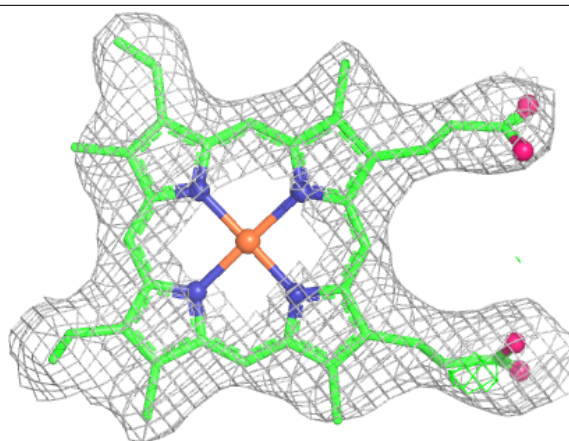
Electron density around HEC B 302:

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



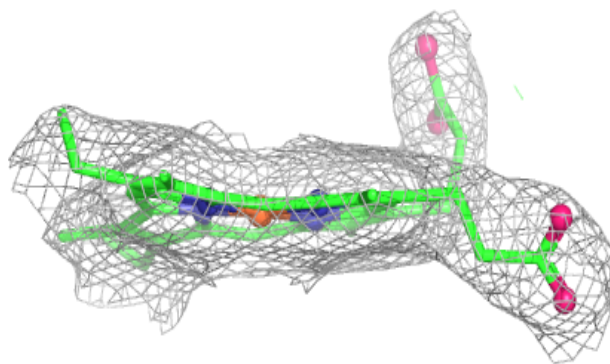
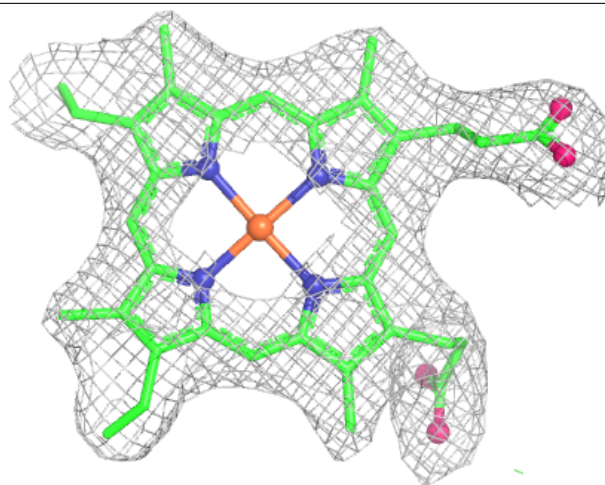
Electron density around HEC C 301:

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



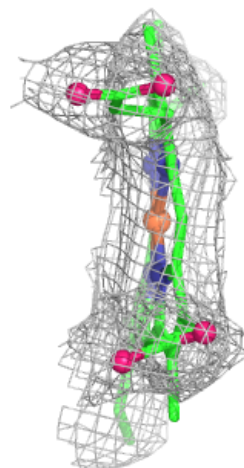
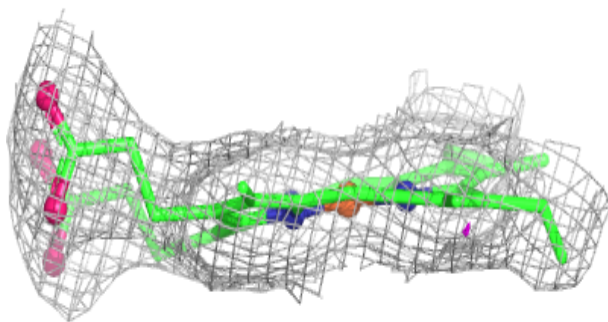
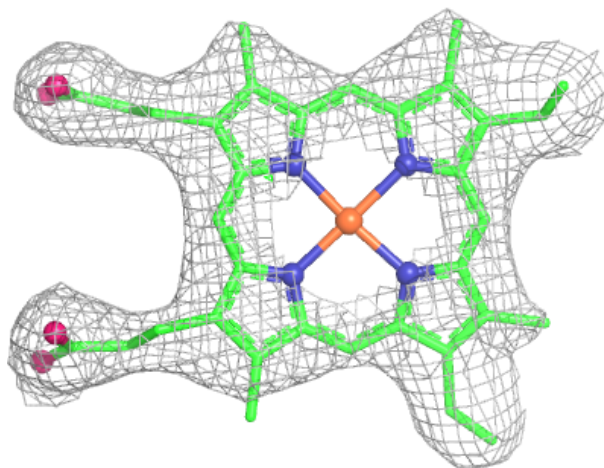
Electron density around HEC C 302:

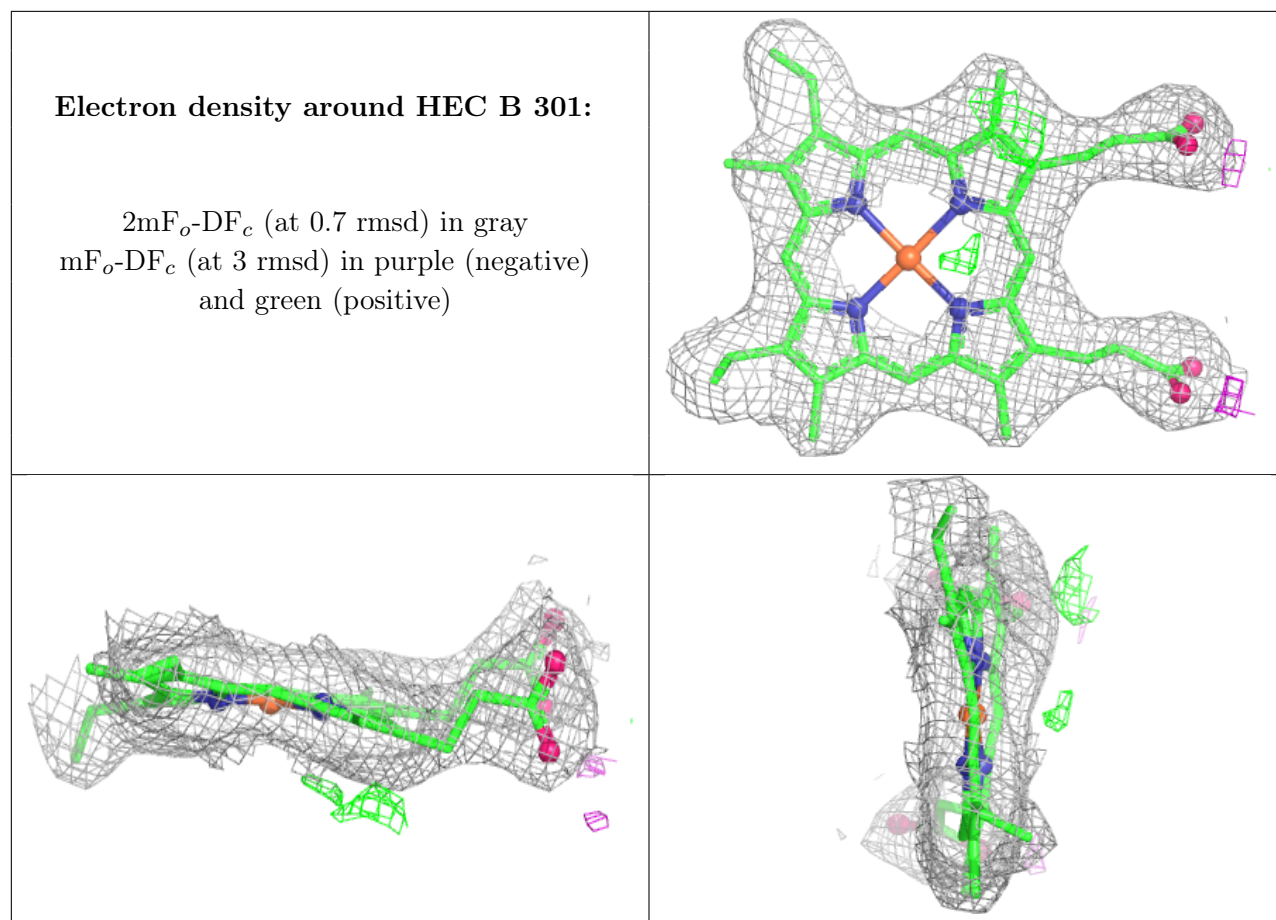
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around HEC A 301:

$2mF_o-DF_c$ (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.