wwPDB X-ray Structure Validation Summary Report (i)

Dec 15, 2024 – 08:37 AM EST

PDB ID	:	7RWY
Title	:	Crystal structure of a Fe-bound RIDC1 variant in the presence of reductant
Authors	:	Kakkis, A.; Golub, E.
Deposited on	:	2021-08-20
Resolution	:	2.20 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 (i) 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 (i)) were used in the production of this report:

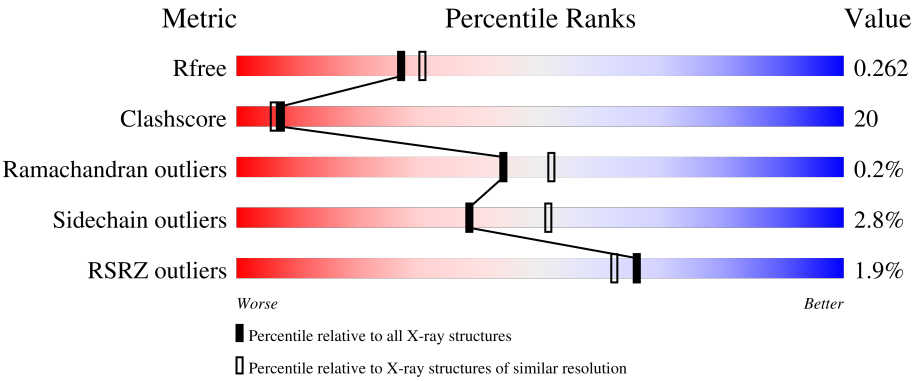
MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.21
EDS	:	3.0
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.004 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.40

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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}	164625	5791 (2.20-2.20)
Clashscore	180529	6634 (2.20-2.20)
Ramachandran outliers	177936	6560 (2.20-2.20)
Sidechain outliers	177891	6561 (2.20-2.20)
RSRZ outliers	164620	5791 (2.20-2.20)

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 <=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	106	2% 81% 17%
1	B	106	% 69% 30%
1	C	106	73% 27%
1	D	106	5% 70% 25% 5%
1	E	106	% 75% 24%

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Mol	Chain	Length	Quality of chain
1	F	106	3% 58% 42%

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	TRS	C	204	-	X	-	-

2 Entry composition (i)

There are 6 unique types of molecules in this entry. The entry contains 5699 atoms, of which 236 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 Soluble cytochrome b562.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	106	Total	C	N	O	S	0	0	0
			815	504	145	160	6			
1	B	106	Total	C	N	O	S	0	0	0
			815	504	145	160	6			
1	C	106	Total	C	N	O	S	0	1	0
			823	510	146	161	6			
1	D	106	Total	C	N	O	S	0	2	0
			832	513	147	166	6			
1	E	106	Total	C	N	O	S	0	1	0
			824	509	146	163	6			
1	F	106	Total	C	N	O	S	0	0	0
			815	504	145	160	6			

There are 78 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	34	ALA	ARG	engineered mutation	UNP P0ABE7
A	38	ALA	LEU	engineered mutation	UNP P0ABE7
A	41	TRP	GLN	engineered mutation	UNP P0ABE7
A	42	SER	LYS	engineered mutation	UNP P0ABE7
A	59	HIS	LYS	engineered mutation	UNP P0ABE7
A	66	TRP	ASP	engineered mutation	UNP P0ABE7
A	69	ILE	VAL	engineered mutation	UNP P0ABE7
A	73	HIS	ASP	engineered mutation	UNP P0ABE7
A	74	ALA	ASP	engineered mutation	UNP P0ABE7
A	77	HIS	LYS	engineered mutation	UNP P0ABE7
A	96	CYS	THR	engineered mutation	UNP P0ABE7
A	98	CYS	ARG	engineered mutation	UNP P0ABE7
A	101	CYS	TYR	engineered mutation	UNP P0ABE7
B	34	ALA	ARG	engineered mutation	UNP P0ABE7
B	38	ALA	LEU	engineered mutation	UNP P0ABE7
B	41	TRP	GLN	engineered mutation	UNP P0ABE7
B	42	SER	LYS	engineered mutation	UNP P0ABE7

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Chain	Residue	Modelled	Actual	Comment	Reference
B	59	HIS	LYS	engineered mutation	UNP P0ABE7
B	66	TRP	ASP	engineered mutation	UNP P0ABE7
B	69	ILE	VAL	engineered mutation	UNP P0ABE7
B	73	HIS	ASP	engineered mutation	UNP P0ABE7
B	74	ALA	ASP	engineered mutation	UNP P0ABE7
B	77	HIS	LYS	engineered mutation	UNP P0ABE7
B	96	CYS	THR	engineered mutation	UNP P0ABE7
B	98	CYS	ARG	engineered mutation	UNP P0ABE7
B	101	CYS	TYR	engineered mutation	UNP P0ABE7
C	34	ALA	ARG	engineered mutation	UNP P0ABE7
C	38	ALA	LEU	engineered mutation	UNP P0ABE7
C	41	TRP	GLN	engineered mutation	UNP P0ABE7
C	42	SER	LYS	engineered mutation	UNP P0ABE7
C	59	HIS	LYS	engineered mutation	UNP P0ABE7
C	66	TRP	ASP	engineered mutation	UNP P0ABE7
C	69	ILE	VAL	engineered mutation	UNP P0ABE7
C	73	HIS	ASP	engineered mutation	UNP P0ABE7
C	74	ALA	ASP	engineered mutation	UNP P0ABE7
C	77	HIS	LYS	engineered mutation	UNP P0ABE7
C	96	CYS	THR	engineered mutation	UNP P0ABE7
C	98	CYS	ARG	engineered mutation	UNP P0ABE7
C	101	CYS	TYR	engineered mutation	UNP P0ABE7
D	34	ALA	ARG	engineered mutation	UNP P0ABE7
D	38	ALA	LEU	engineered mutation	UNP P0ABE7
D	41	TRP	GLN	engineered mutation	UNP P0ABE7
D	42	SER	LYS	engineered mutation	UNP P0ABE7
D	59	HIS	LYS	engineered mutation	UNP P0ABE7
D	66	TRP	ASP	engineered mutation	UNP P0ABE7
D	69	ILE	VAL	engineered mutation	UNP P0ABE7
D	73	HIS	ASP	engineered mutation	UNP P0ABE7
D	74	ALA	ASP	engineered mutation	UNP P0ABE7
D	77	HIS	LYS	engineered mutation	UNP P0ABE7
D	96	CYS	THR	engineered mutation	UNP P0ABE7
D	98	CYS	ARG	engineered mutation	UNP P0ABE7
D	101	CYS	TYR	engineered mutation	UNP P0ABE7
E	34	ALA	ARG	engineered mutation	UNP P0ABE7
E	38	ALA	LEU	engineered mutation	UNP P0ABE7
E	41	TRP	GLN	engineered mutation	UNP P0ABE7
E	42	SER	LYS	engineered mutation	UNP P0ABE7
E	59	HIS	LYS	engineered mutation	UNP P0ABE7
E	66	TRP	ASP	engineered mutation	UNP P0ABE7
E	69	ILE	VAL	engineered mutation	UNP P0ABE7

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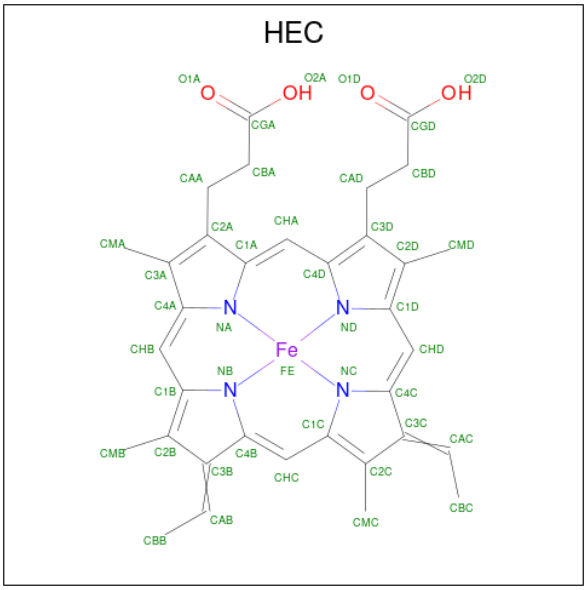
Chain	Residue	Modelled	Actual	Comment	Reference
E	73	HIS	ASP	engineered mutation	UNP P0ABE7
E	74	ALA	ASP	engineered mutation	UNP P0ABE7
E	77	HIS	LYS	engineered mutation	UNP P0ABE7
E	96	CYS	THR	engineered mutation	UNP P0ABE7
E	98	CYS	ARG	engineered mutation	UNP P0ABE7
E	101	CYS	TYR	engineered mutation	UNP P0ABE7
F	34	ALA	ARG	engineered mutation	UNP P0ABE7
F	38	ALA	LEU	engineered mutation	UNP P0ABE7
F	41	TRP	GLN	engineered mutation	UNP P0ABE7
F	42	SER	LYS	engineered mutation	UNP P0ABE7
F	59	HIS	LYS	engineered mutation	UNP P0ABE7
F	66	TRP	ASP	engineered mutation	UNP P0ABE7
F	69	ILE	VAL	engineered mutation	UNP P0ABE7
F	73	HIS	ASP	engineered mutation	UNP P0ABE7
F	74	ALA	ASP	engineered mutation	UNP P0ABE7
F	77	HIS	LYS	engineered mutation	UNP P0ABE7
F	96	CYS	THR	engineered mutation	UNP P0ABE7
F	98	CYS	ARG	engineered mutation	UNP P0ABE7
F	101	CYS	TYR	engineered mutation	UNP P0ABE7

- Molecule 2 is HEME C (three-letter code: HEC) (formula: $C_{34}H_{34}FeN_4O_4$).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total	C	Fe	H	N	O	0	0
			73	34	1	30	4	4		
2	B	1	Total	C	Fe	H	N	O	0	0
			73	34	1	30	4	4		

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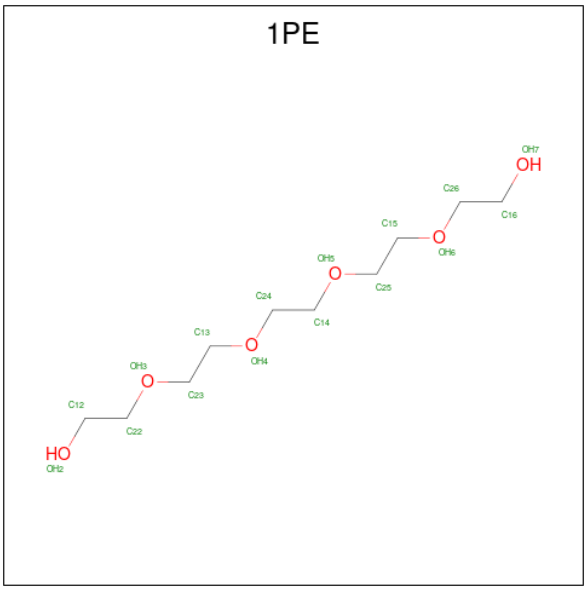
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	C	1	Total	C	Fe	H	N	O	0	0
			73	34	1	30	4	4		
2	D	1	Total	C	Fe	H	N	O	0	0
			73	34	1	30	4	4		
2	E	1	Total	C	Fe	H	N	O	0	0
			73	34	1	30	4	4		
2	F	1	Total	C	Fe	H	N	O	0	0
			73	34	1	30	4	4		

- Molecule 3 is PENTAETHYLENE GLYCOL (three-letter code: 1PE) (formula: $C_{10}H_{22}O_6$).

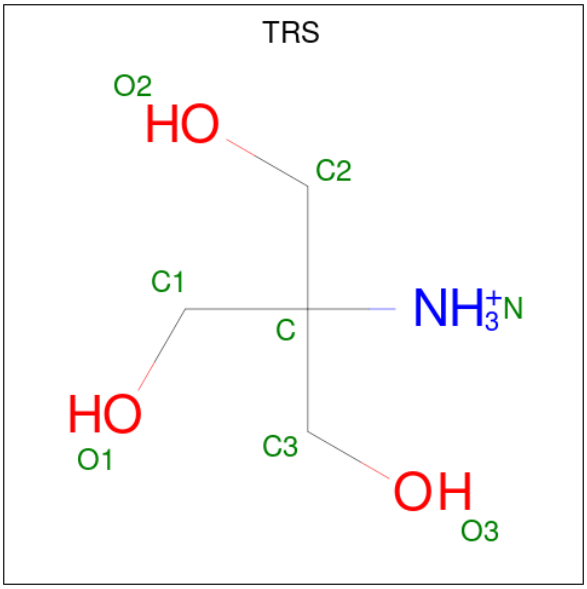


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	B	1	Total	C	H	O	0	0
			38	10	22	6		
3	D	1	Total	C	H	O	0	0
			38	10	22	6		

- Molecule 4 is FE (III) ION (three-letter code: FE) (formula: Fe) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	C	2	Total	Fe	0	0
			2	2		
4	D	2	Total	Fe	0	0
			2	2		

- Molecule 5 is 2-AMINO-2-HYDROXYMETHYL-PROPANE-1,3-DIOL (three-letter code: TRS) (formula: $C_4H_{12}NO_3$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	C	1	Total	C	H	N	O	0	0
			20	4	12	1	3		

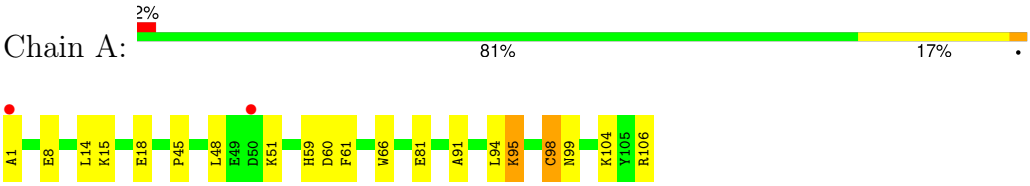
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	31	Total	O	0	0
			31	31		
6	B	47	Total	O	0	0
			47	47		
6	C	51	Total	O	0	0
			51	51		
6	D	39	Total	O	0	0
			39	39		
6	E	45	Total	O	0	0
			45	45		
6	F	24	Total	O	0	0
			24	24		

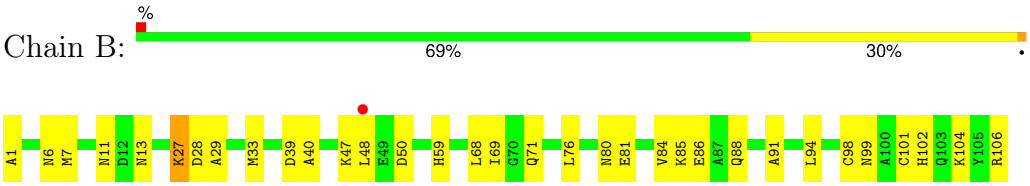
3 Residue-property plots (i)

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.

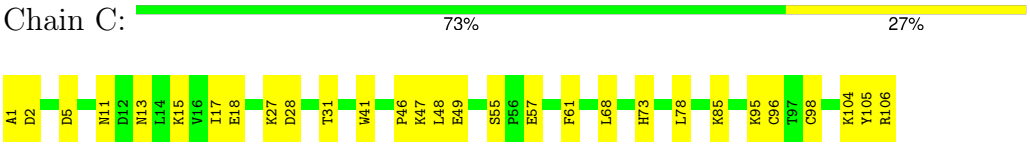
• Molecule 1: Soluble cytochrome b562



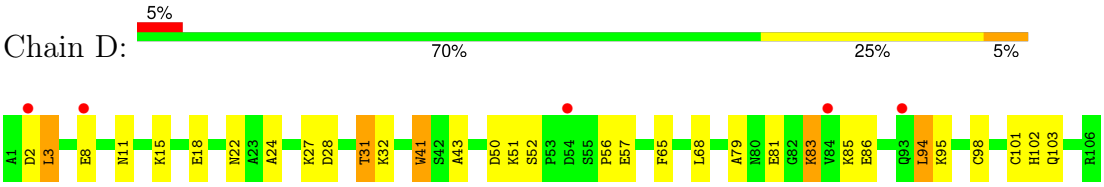
• Molecule 1: Soluble cytochrome b562



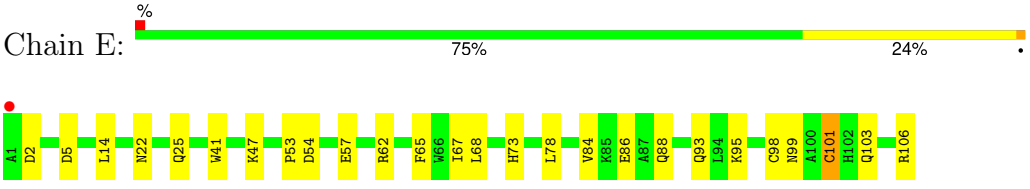
• Molecule 1: Soluble cytochrome b562



• Molecule 1: Soluble cytochrome b562

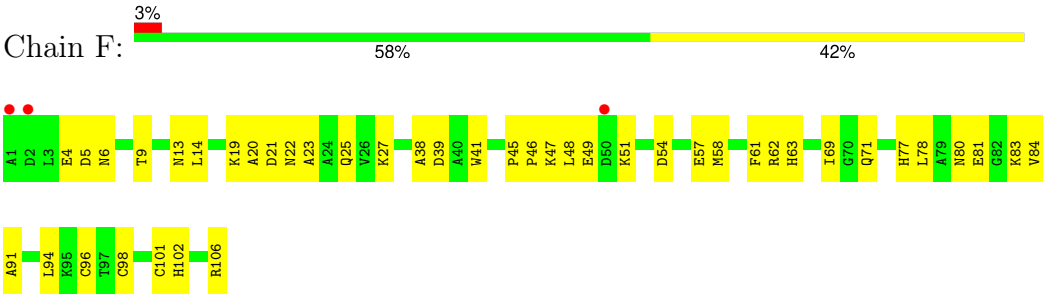


• Molecule 1: Soluble cytochrome b562



• Molecule 1: Soluble cytochrome b562

Chain F:


3%
58%
42%
A1
D2
L3
E4
D5
N6
T9
N13
L14
K19
A20
D21
N22
A23
A24
Q25
V26
K27
A38
D39
A40
W41
P45
P46
K47
L48
E49
D50
K51
D54
E57
M58
F61
R62
H63
I69
G70
Q71
H77
L78
A79
N80
E81
G82
K83
V84
K85
Q88
A91
L94
K95
C96
T97
C98
C101
H102
R106

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	52.55Å 82.68Å 74.14Å 90.00° 95.22° 90.00°	Depositor
Resolution (Å)	44.65 – 2.20 44.65 – 2.20	Depositor EDS
% Data completeness (in resolution range)	94.4 (44.65-2.20) 93.7 (44.65-2.20)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	4.26 (at 2.20Å)	Xtriage
Refinement program	PHENIX 1.18.2_3874	Depositor
R, R_{free}	0.202 , 0.259 0.206 , 0.262	Depositor DCC
R_{free} test set	1598 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å^2)	31.0	Xtriage
Anisotropy	0.409	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 42.7	EDS
L-test for twinning[2]	$< \|L\| > = 0.47$, $< L^2 > = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	5699	wwPDB-VP
Average B, all atoms (Å^2)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 32.83 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 8.8147e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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: HEC, FE, TRS, 1PE

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	0.37	0/830	0.49	0/1121
1	B	0.40	0/830	0.51	0/1121
1	C	0.42	0/838	0.56	0/1132
1	D	0.41	0/847	0.54	0/1144
1	E	0.45	0/839	0.54	1/1133 (0.1%)
1	F	0.37	0/830	0.52	0/1121
All	All	0.40	0/5014	0.53	1/6772 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	E	101	CYS	CA-CB-SG	-5.96	103.28	114.00

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	815	0	791	23	1
1	B	815	0	792	39	0
1	C	823	0	800	29	0
1	D	832	0	800	40	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	E	824	0	796	28	0
1	F	815	0	791	37	1
2	A	43	30	31	8	0
2	B	43	30	32	15	0
2	C	43	30	30	7	0
2	D	43	30	32	8	0
2	E	43	30	31	10	0
2	F	43	30	31	9	0
3	B	16	22	22	1	0
3	D	16	22	22	1	0
4	C	2	0	0	0	0
4	D	2	0	0	0	0
5	C	8	12	12	4	0
6	A	31	0	0	7	2
6	B	47	0	0	10	1
6	C	51	0	0	10	0
6	D	39	0	0	11	0
6	E	45	0	0	4	1
6	F	24	0	0	5	0
All	All	5463	236	5013	208	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 20.

The worst 5 of 208 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)
5:C:204:TRS:N	5:C:204:TRS:C	1.71	1.47
1:E:78:LEU:HD13	1:E:86[A]:GLU:HG3	1.46	0.96
1:D:22:ASN:ND2	6:D:301:HOH:O	1.98	0.94
1:B:101:CYS:SG	2:B:202:HEC:CAC	2.62	0.87
1:F:96:CYS:SG	6:F:323:HOH:O	2.32	0.87

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:A:81:GLU:OE1	1:F:47:LYS:NZ[2_854]	2.03	0.17
6:A:308:HOH:O	6:B:306:HOH:O[2_854]	2.12	0.08
6:A:330:HOH:O	6:E:336:HOH:O[1_655]	2.15	0.05

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	104/106 (98%)	102 (98%)	2 (2%)	0	100 100
1	B	104/106 (98%)	101 (97%)	3 (3%)	0	100 100
1	C	105/106 (99%)	105 (100%)	0	0	100 100
1	D	106/106 (100%)	100 (94%)	5 (5%)	1 (1%)	14 14
1	E	105/106 (99%)	104 (99%)	1 (1%)	0	100 100
1	F	104/106 (98%)	100 (96%)	4 (4%)	0	100 100
All	All	628/636 (99%)	612 (98%)	15 (2%)	1 (0%)	44 52

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	83	LYS

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	83/83 (100%)	81 (98%)	2 (2%)	44 57
1	B	83/83 (100%)	81 (98%)	2 (2%)	44 57
1	C	84/83 (101%)	82 (98%)	2 (2%)	44 57
1	D	85/83 (102%)	79 (93%)	6 (7%)	12 13
1	E	84/83 (101%)	83 (99%)	1 (1%)	67 80
1	F	83/83 (100%)	82 (99%)	1 (1%)	67 80

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	502/498 (101%)	488 (97%)	14 (3%)	38 51

5 of 14 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	3	LEU
1	D	31	THR
1	F	21	ASP
1	D	94	LEU
1	E	103	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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)

Of 13 ligands modelled in this entry, 4 are monoatomic - leaving 9 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	202	1	32,50,50	2.19	5 (15%)	30,82,82	1.58	6 (20%)
2	HEC	C	203	1	32,50,50	2.30	4 (12%)	30,82,82	1.68	8 (26%)
2	HEC	E	201	1	32,50,50	2.04	4 (12%)	30,82,82	1.81	6 (20%)
2	HEC	A	201	1	32,50,50	2.22	4 (12%)	30,82,82	1.76	7 (23%)
3	1PE	D	203	-	15,15,15	0.55	0	14,14,14	0.46	0
2	HEC	F	201	1	32,50,50	2.06	4 (12%)	30,82,82	1.55	5 (16%)
3	1PE	B	201	-	15,15,15	0.50	0	14,14,14	0.43	0
5	TRS	C	204	-	7,7,7	3.35	4 (57%)	9,9,9	5.82	7 (77%)
2	HEC	D	204	1	32,50,50	2.04	5 (15%)	30,82,82	1.80	9 (30%)

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	202	1	-	2/10/54/54	-
2	HEC	C	203	1	-	0/10/54/54	-
2	HEC	E	201	1	-	0/10/54/54	-
2	HEC	A	201	1	-	0/10/54/54	-
3	1PE	D	203	-	-	9/13/13/13	-
2	HEC	F	201	1	-	3/10/54/54	-
3	1PE	B	201	-	-	6/13/13/13	-
5	TRS	C	204	-	-	5/9/9/9	-
2	HEC	D	204	1	-	1/10/54/54	-

The worst 5 of 30 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	201	HEC	C3C-C2C	-7.21	1.32	1.40
5	C	204	TRS	C-N	6.96	1.71	1.49
2	C	203	HEC	C2B-C3B	-6.91	1.33	1.40
2	B	202	HEC	C2B-C3B	-6.36	1.33	1.40
2	D	204	HEC	C2B-C3B	-6.29	1.33	1.40

The worst 5 of 48 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
5	C	204	TRS	C3-C-N	9.70	132.92	108.17

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
5	C	204	TRS	C2-C-C1	-7.13	91.66	110.66
5	C	204	TRS	C3-C-C1	-6.84	92.43	110.66
5	C	204	TRS	C2-C-N	6.80	125.50	108.17
5	C	204	TRS	C3-C-C2	-6.27	93.96	110.66

There are no chirality outliers.

5 of 26 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	C	204	TRS	C1-C-C2-O2
3	B	201	1PE	OH6-C15-C25-OH5
3	D	203	1PE	OH4-C13-C23-OH3
3	B	201	1PE	C13-C23-OH3-C22
3	B	201	1PE	OH4-C13-C23-OH3

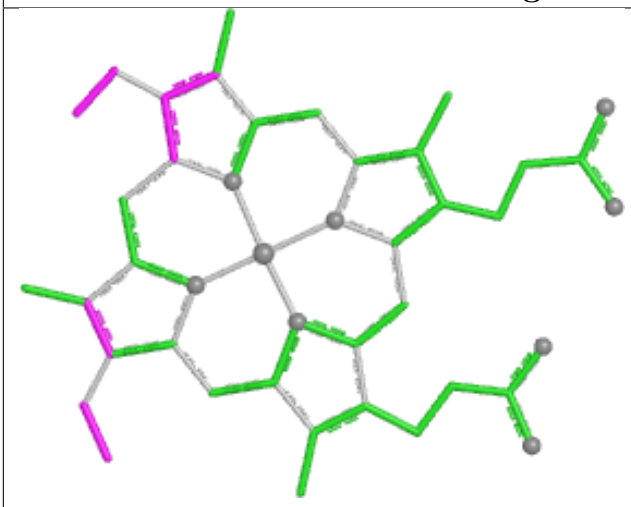
There are no ring outliers.

9 monomers are involved in 63 short contacts:

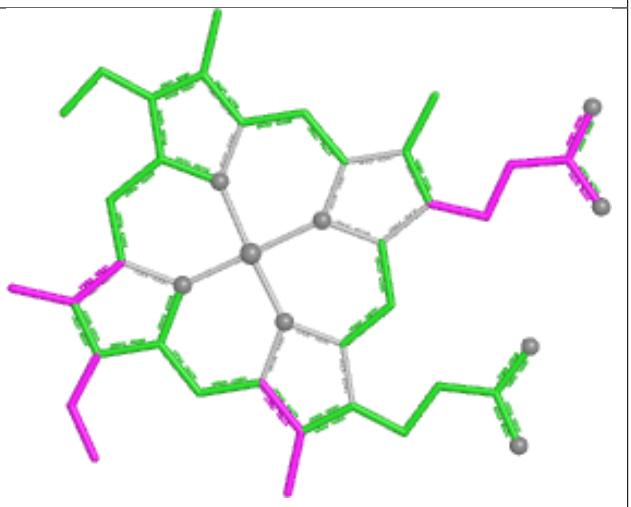
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	202	HEC	15	0
2	C	203	HEC	7	0
2	E	201	HEC	10	0
2	A	201	HEC	8	0
3	D	203	1PE	1	0
2	F	201	HEC	9	0
3	B	201	1PE	1	0
5	C	204	TRS	4	0
2	D	204	HEC	8	0

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 then 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.

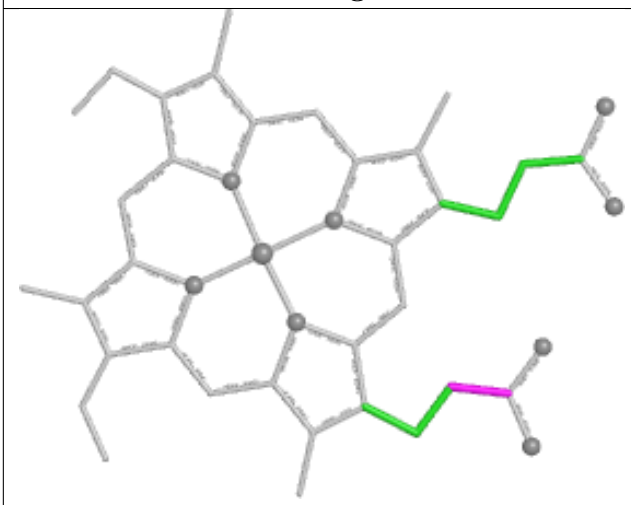
Ligand HEC B 202



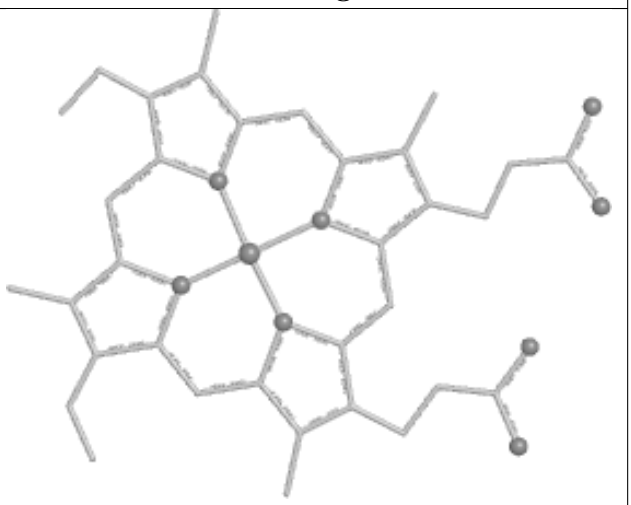
Bond lengths



Bond angles

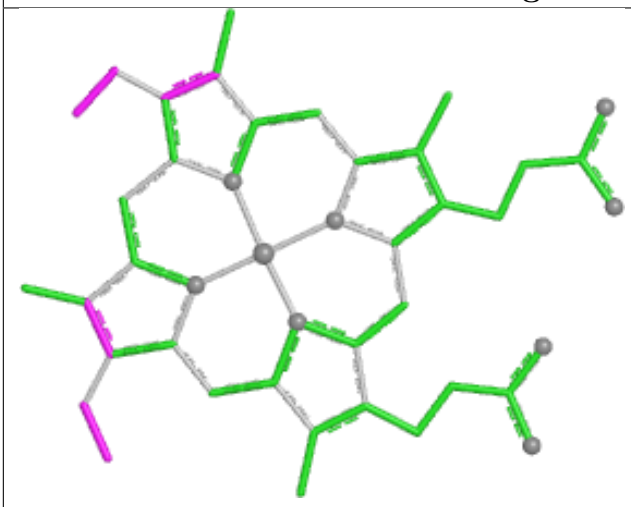


Torsions



Rings

Ligand HEC C 203



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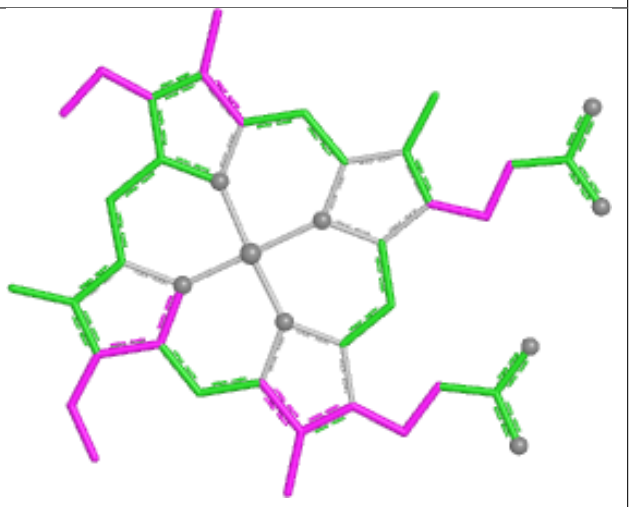
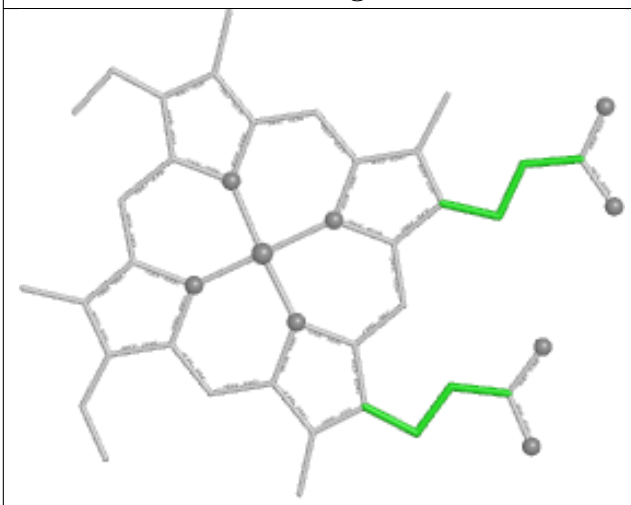
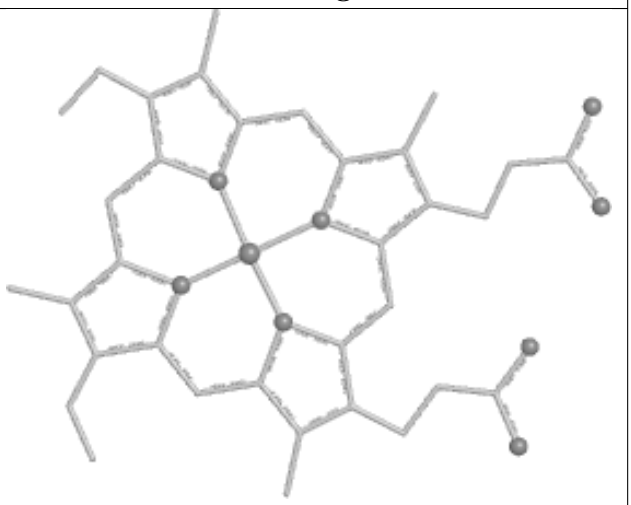


Bond angles

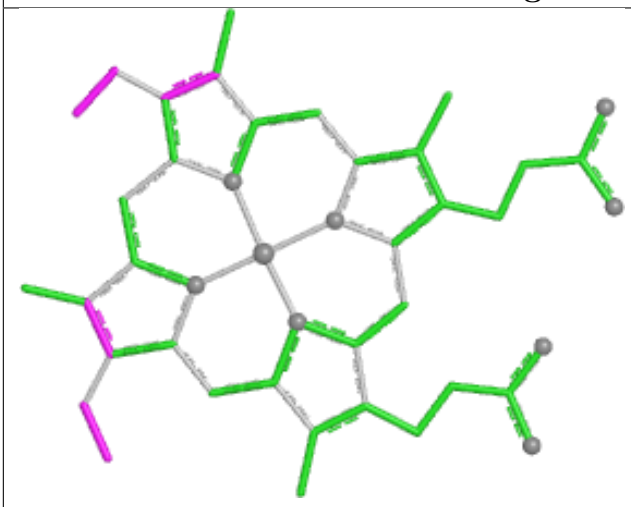


Torsions

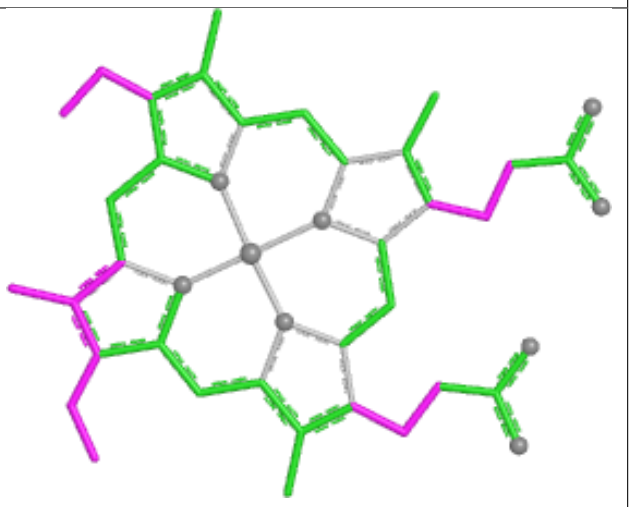


Rings

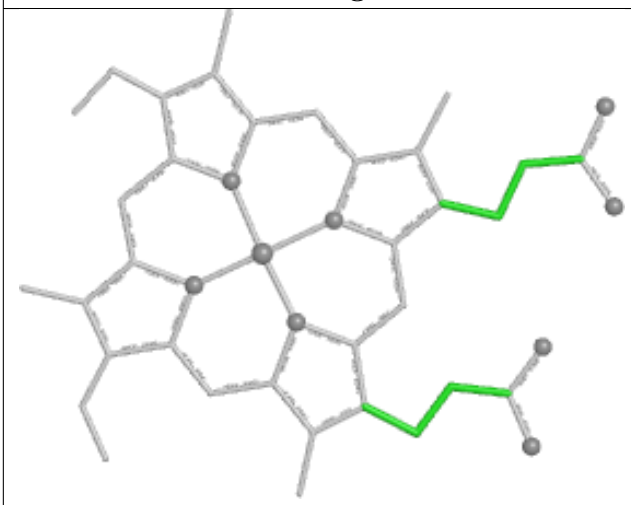
Ligand HEC E 201



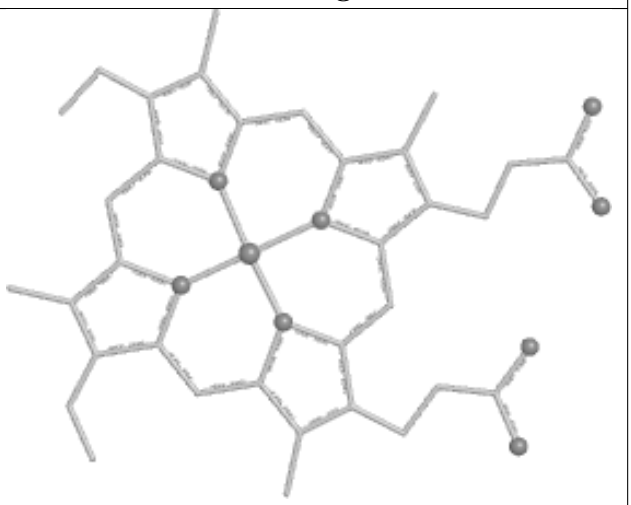
Bond lengths



Bond angles

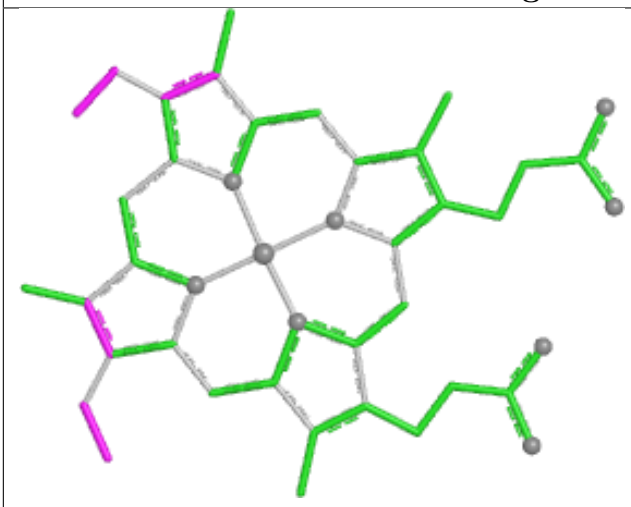


Torsions

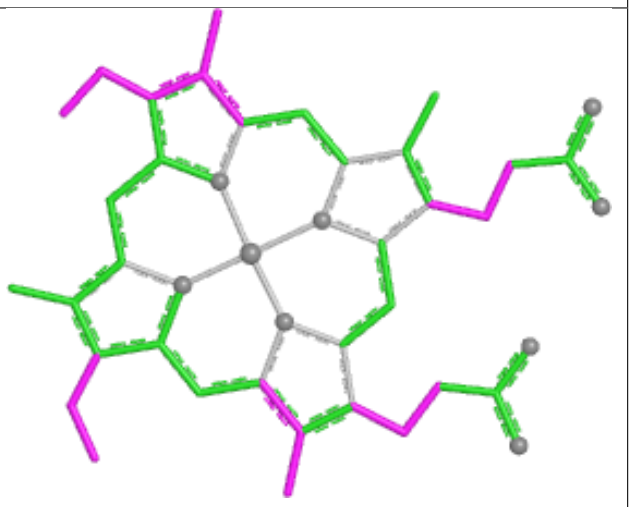


Rings

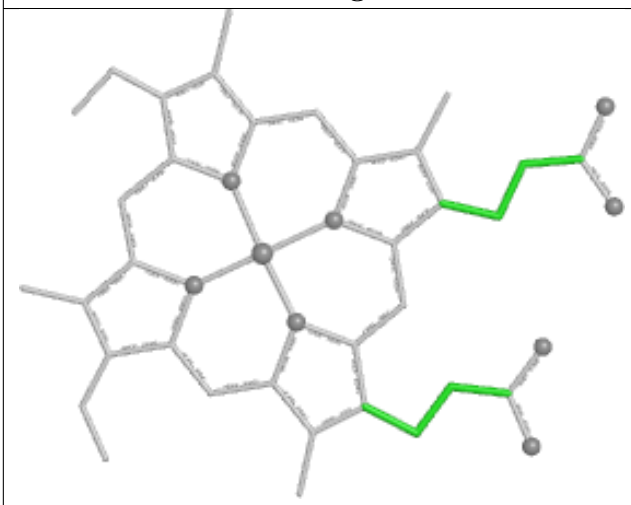
Ligand HEC A 201



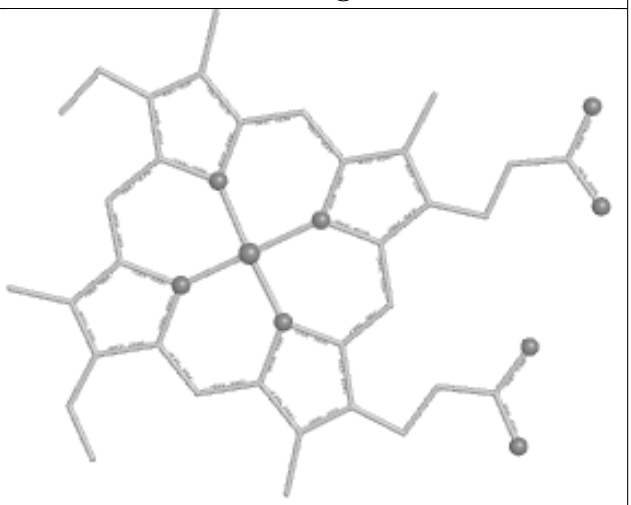
Bond lengths



Bond angles

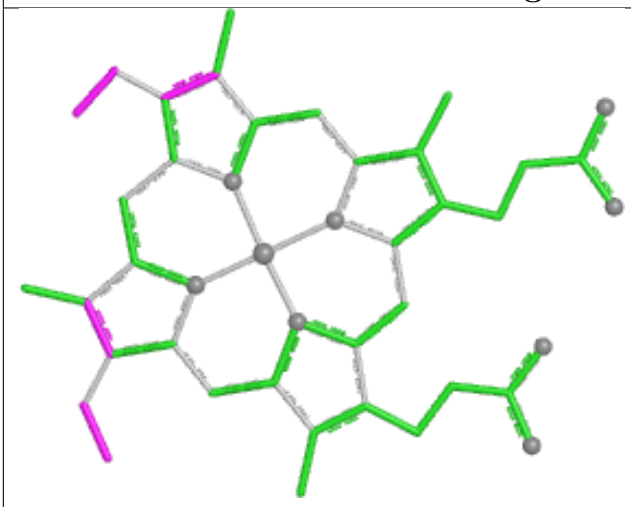


Torsions

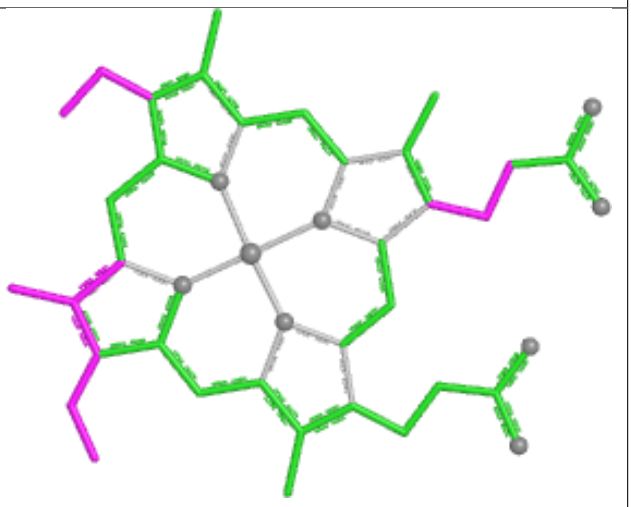


Rings

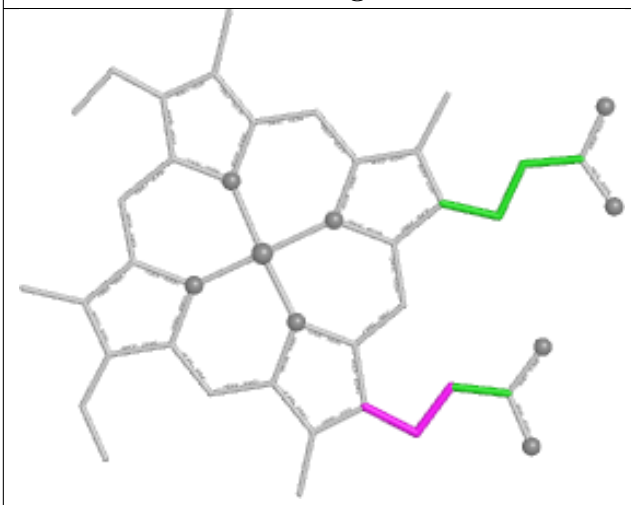
Ligand HEC F 201



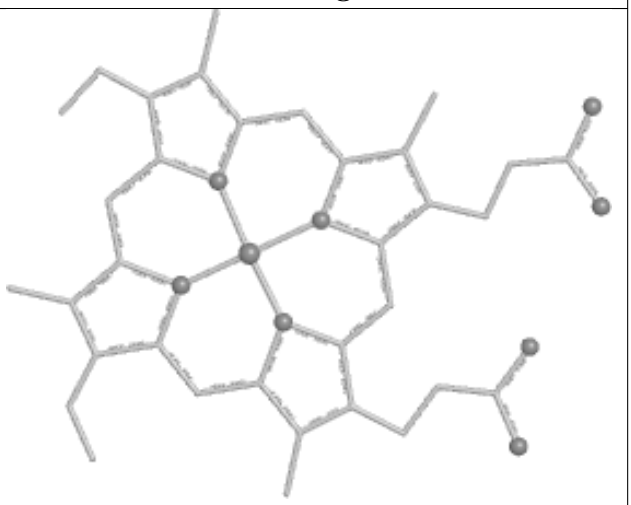
Bond lengths



Bond angles

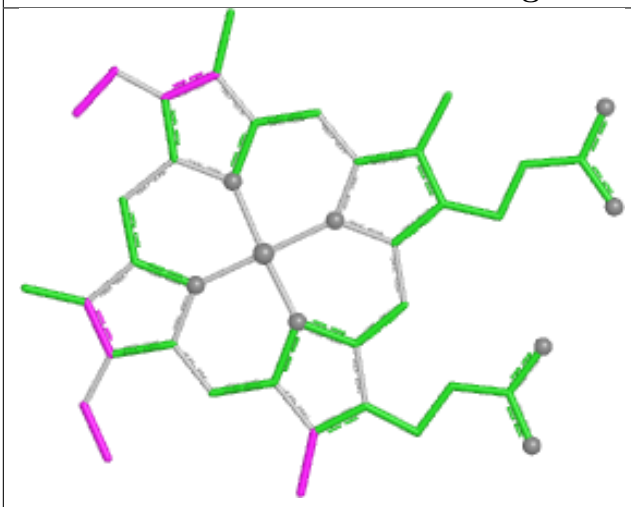


Torsions

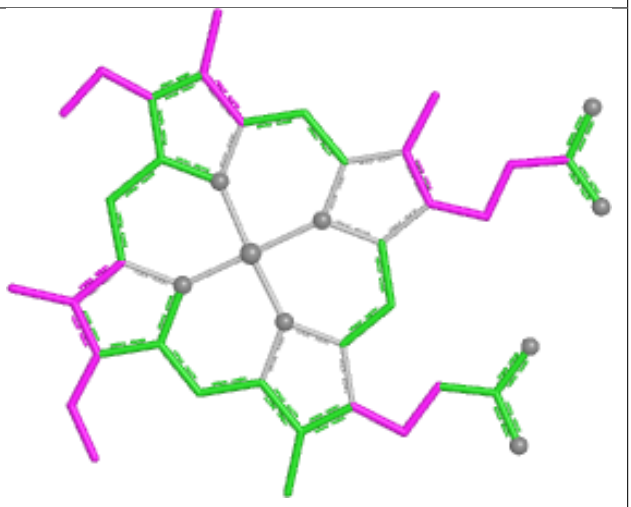


Rings

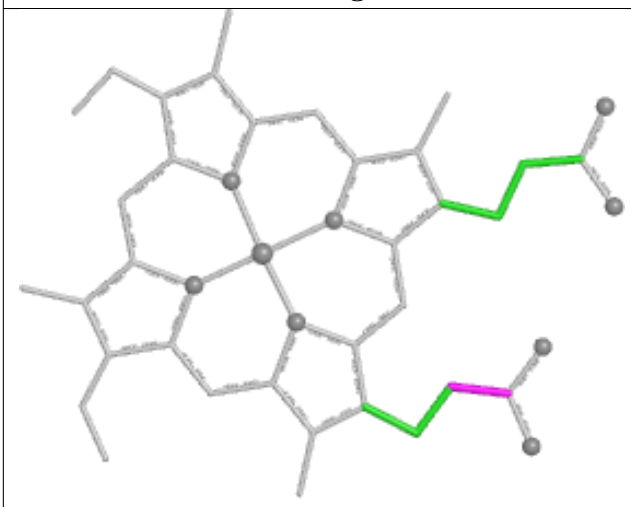
Ligand HEC D 204



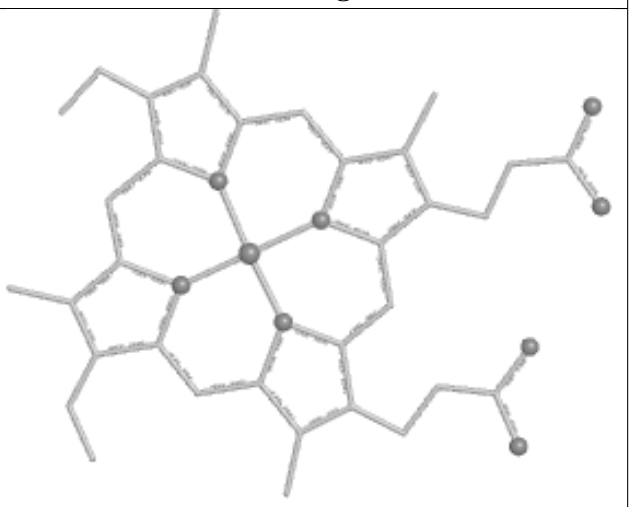
Bond lengths



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, 95^{th} 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	106/106 (100%)	0.23	2 (1%) 66 62	22, 41, 57, 66	0
1	B	106/106 (100%)	0.22	1 (0%) 81 78	20, 37, 55, 64	0
1	C	106/106 (100%)	-0.04	0 100 100	13, 32, 49, 53	1 (0%)
1	D	106/106 (100%)	0.33	5 (4%) 37 34	20, 39, 58, 73	2 (1%)
1	E	106/106 (100%)	-0.04	1 (0%) 81 78	21, 35, 52, 62	1 (0%)
1	F	106/106 (100%)	0.37	3 (2%) 55 52	18, 45, 68, 84	0
All	All	636/636 (100%)	0.18	12 (1%) 66 62	13, 37, 58, 84	4 (0%)

The worst 5 of 12 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	1	ALA	3.3
1	E	1	ALA	3.1
1	A	50	ASP	2.6
1	A	1	ALA	2.5
1	F	50	ASP	2.5

6.2 Non-standard residues in protein, DNA, RNA chains

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

6.3 Carbohydrates

There are no monosaccharides in this entry.

6.4 Ligands (i)

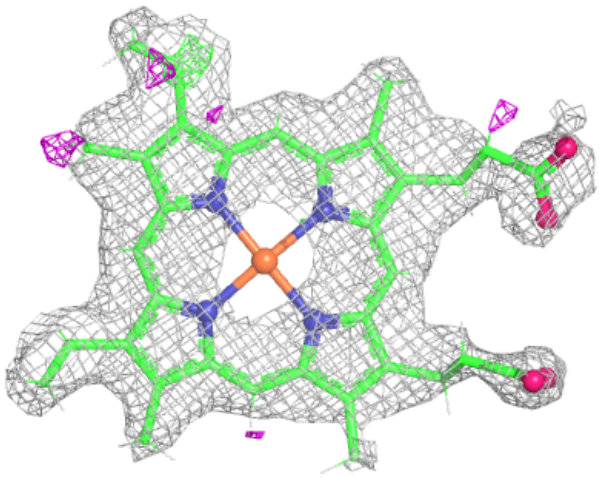
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

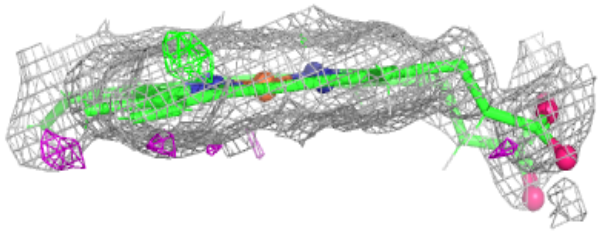
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
5	TRS	C	204	8/8	0.73	0.14	39,64,86,86	0
3	1PE	D	203	16/16	0.85	0.12	31,46,56,58	0
3	1PE	B	201	16/16	0.88	0.10	26,41,53,58	0
2	HEC	F	201	43/43	0.94	0.11	33,47,76,83	0
2	HEC	B	202	43/43	0.95	0.11	30,44,60,70	0
2	HEC	D	204	43/43	0.95	0.09	26,35,51,57	0
2	HEC	A	201	43/43	0.96	0.10	27,39,57,58	0
2	HEC	E	201	43/43	0.96	0.09	21,33,68,89	0
2	HEC	C	203	43/43	0.96	0.08	18,34,46,52	0
4	FE	C	202	1/1	0.98	0.03	32,32,32,32	0
4	FE	C	201	1/1	0.98	0.03	25,25,25,25	0
4	FE	D	202	1/1	0.99	0.03	30,30,30,30	0
4	FE	D	201	1/1	0.99	0.03	27,27,27,27	0

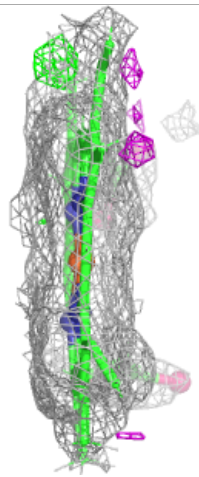
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around HEC F 201:

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

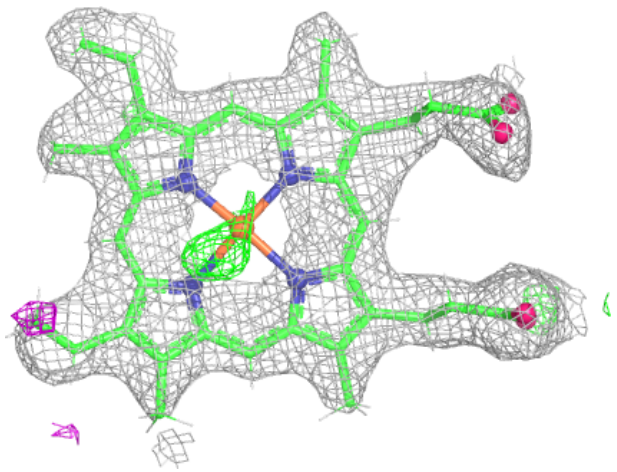


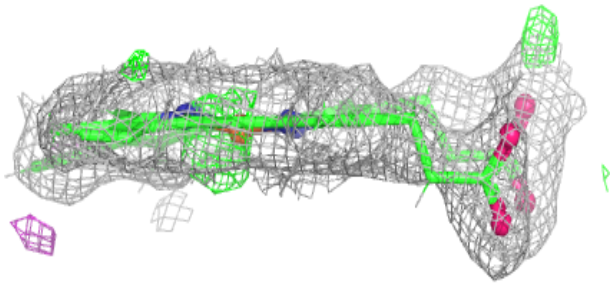


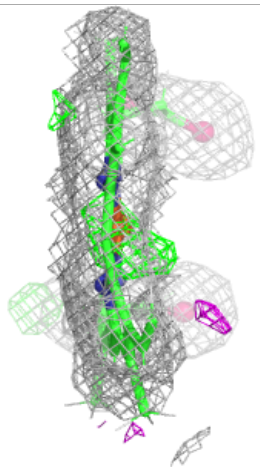


Electron density around HEC B 202:

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

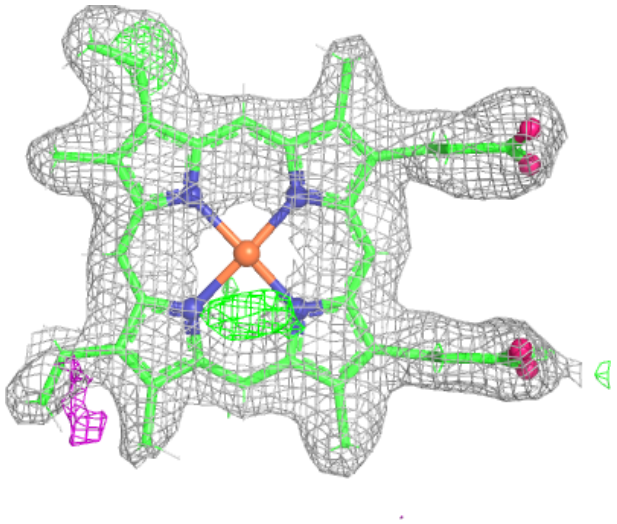


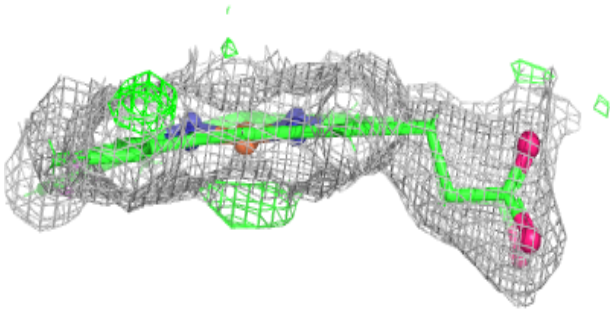


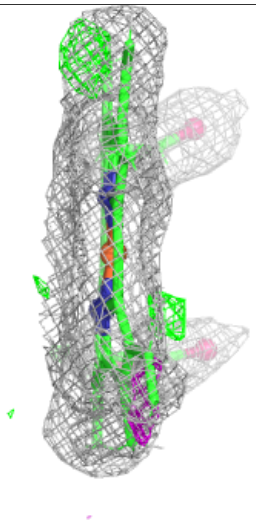


Electron density around HEC D 204:

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

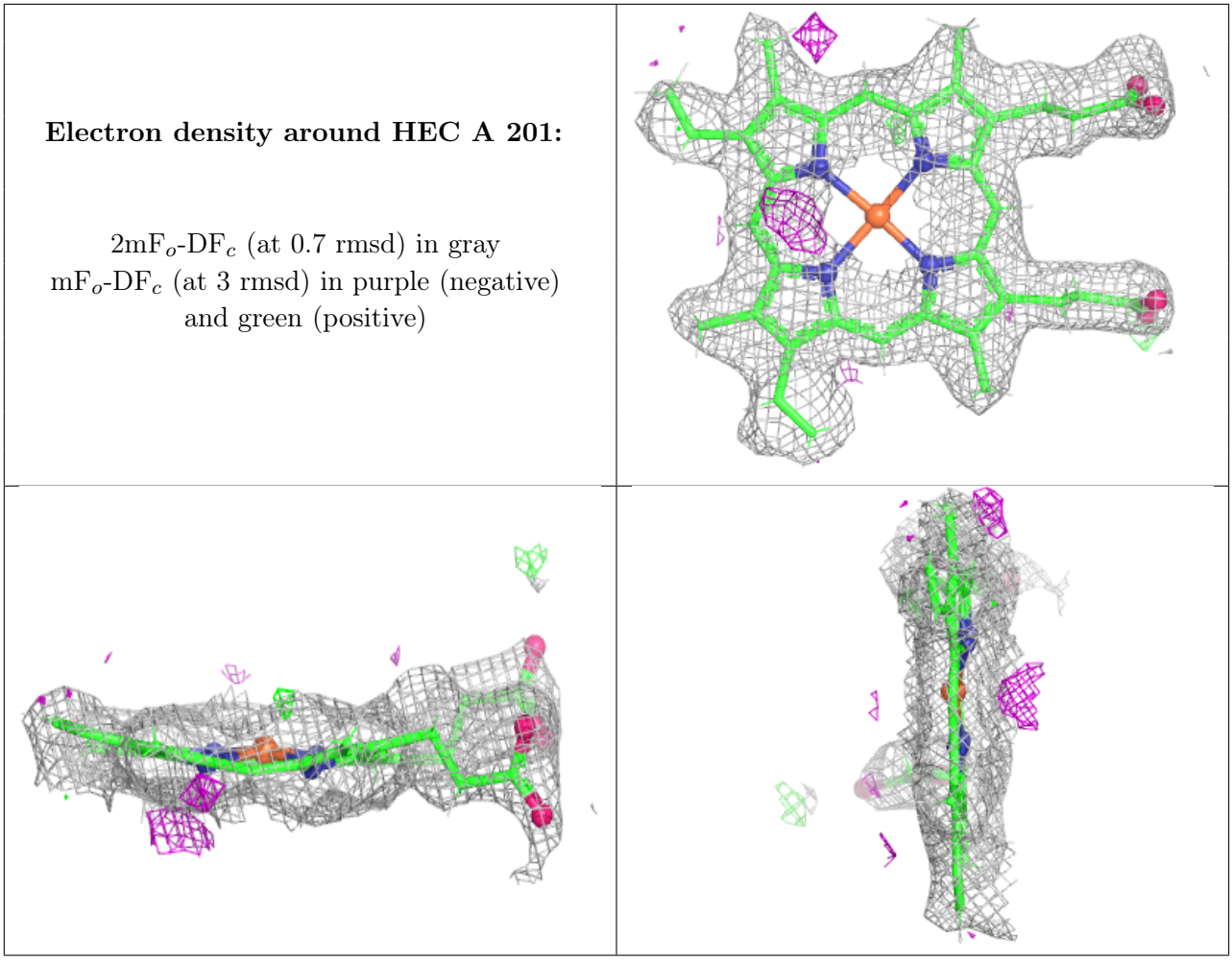






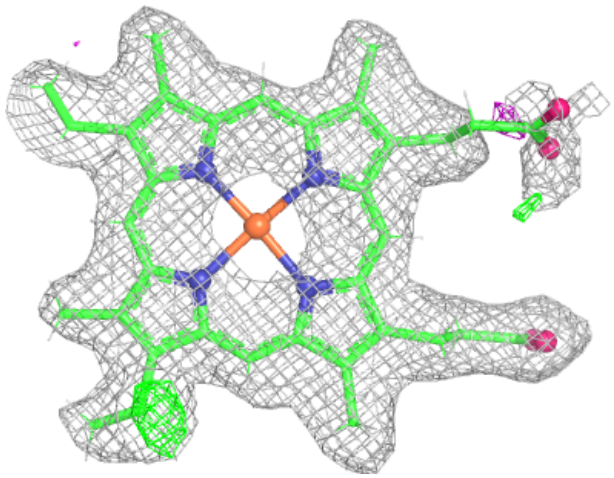
Electron density around HEC A 201:

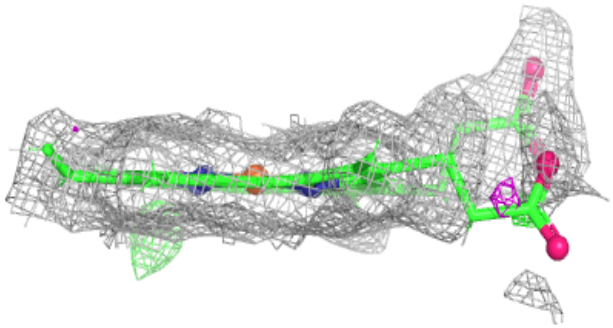
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

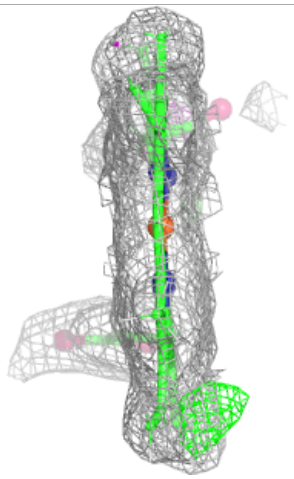


Electron density around HEC E 201:

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

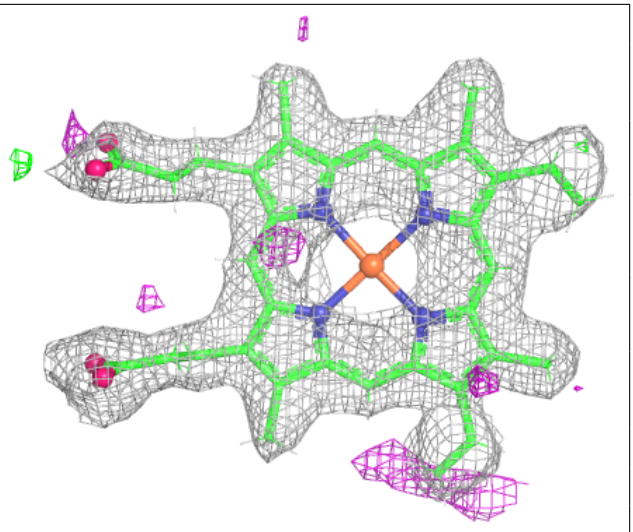


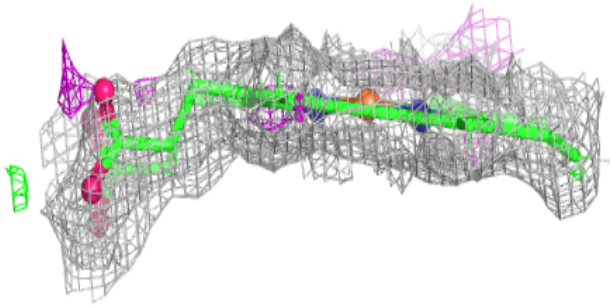


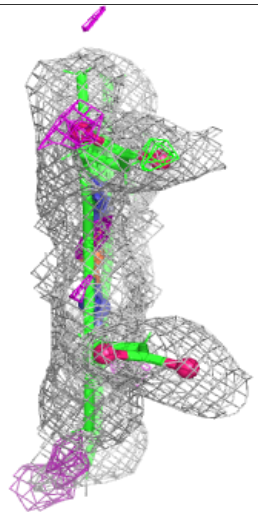


Electron density around HEC C 203:

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

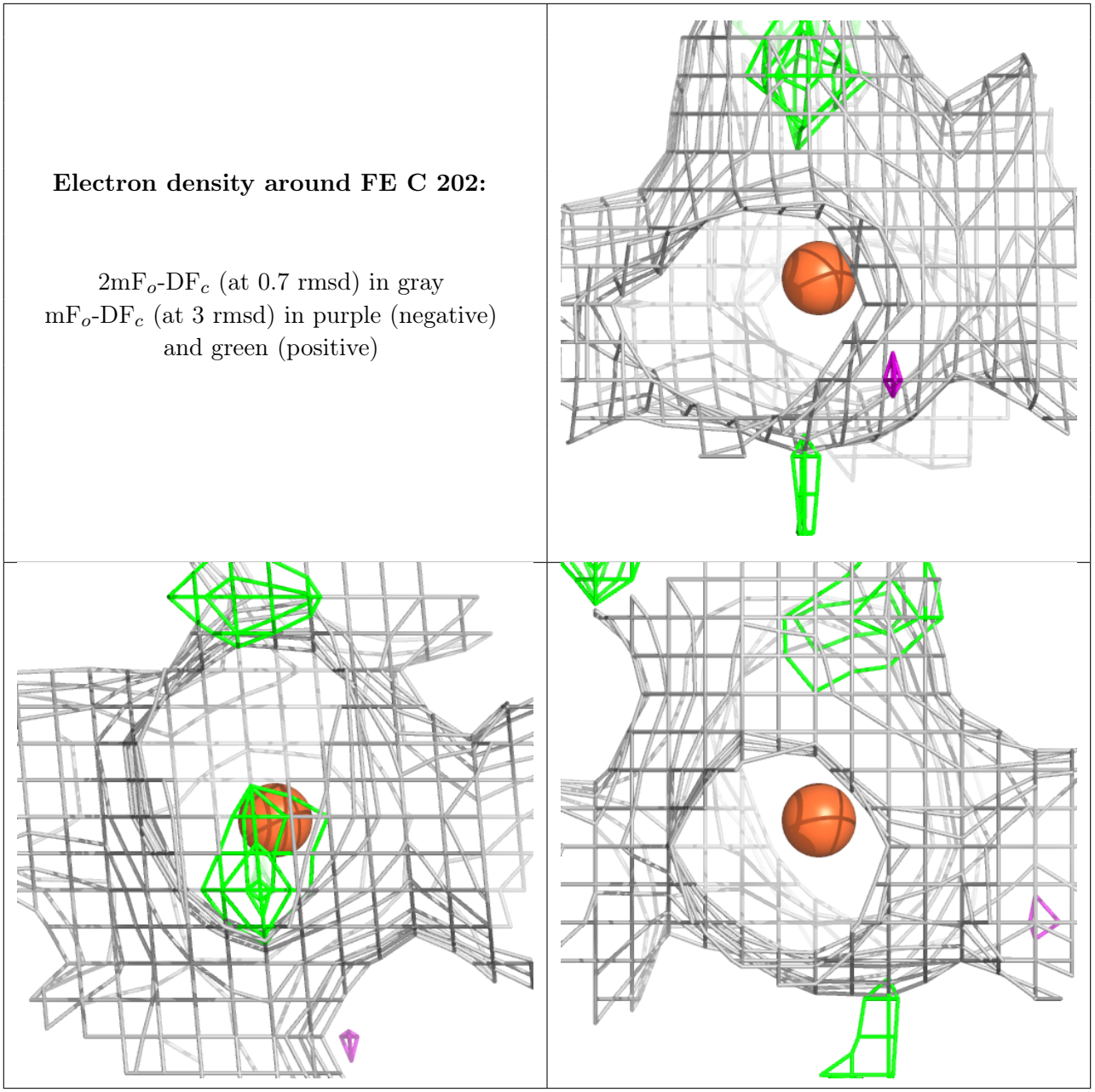






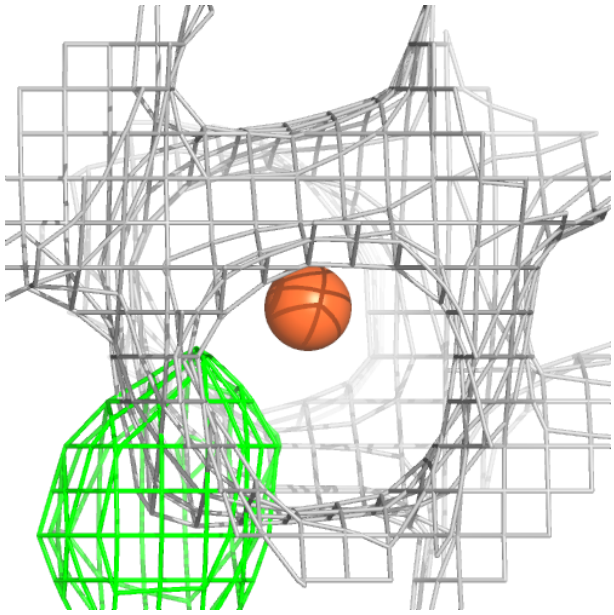
Electron density around FE C 202:

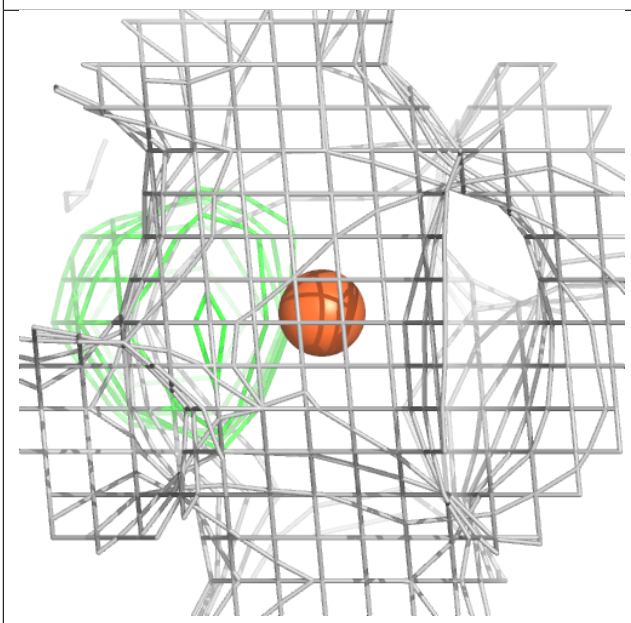
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

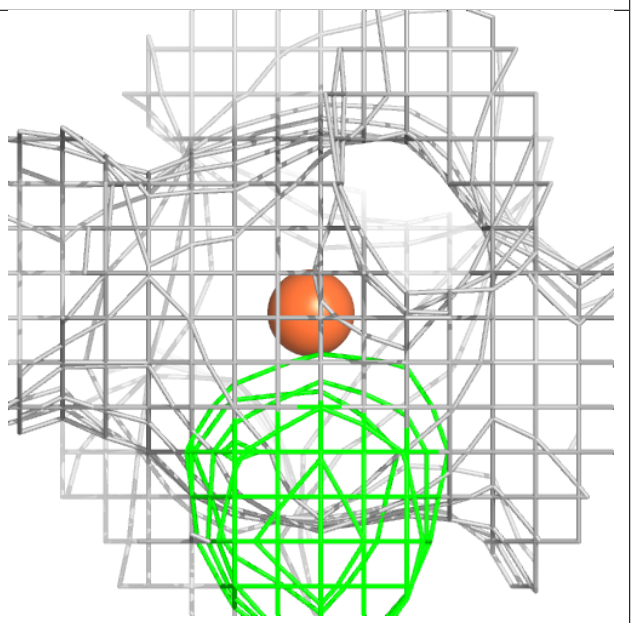


Electron density around FE C 201:

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

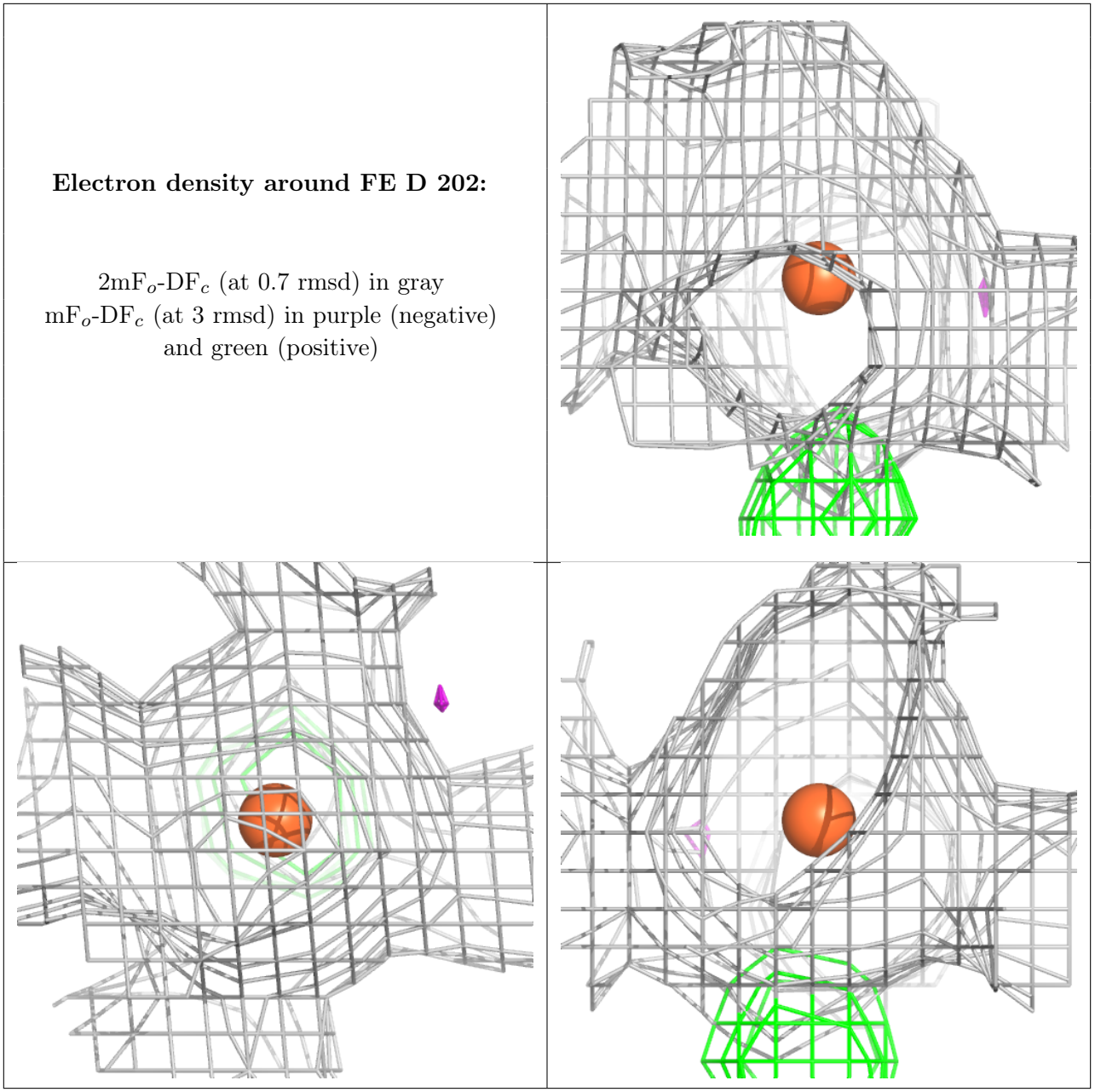






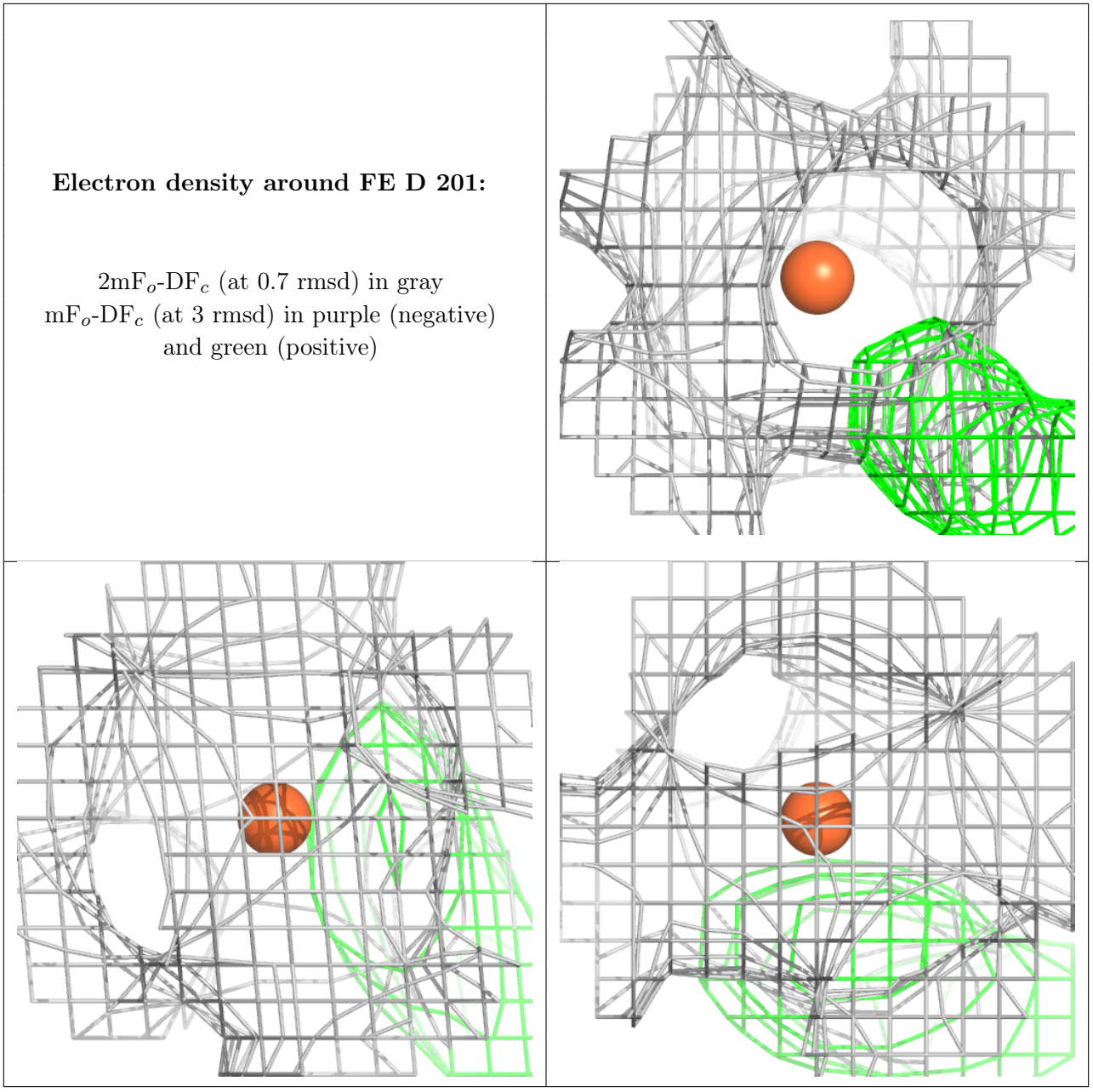
Electron density around FE D 202:

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



Electron density around FE D 201:

$2mF_o$-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.