Full wwPDB X-ray Structure Validation Report (i)

Nov 11, 2023 – 07:16 pm GMT

PDB ID	:	1E39
Title	:	Flavocytochrome C3 from Shewanella frigidimarina histidine 365 mutated to alanine
Authors	:	Doherty, M.K.; Pealing, S.L.; Miles, C.S.; Moysey, R.; Taylor, P.; Walkinshaw, M.D.; Reid, G.A.; Chapman, S.K.
Deposited on	:	2000-06-07
Resolution	:	1.80 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 (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 (1)) were used in the production of this report:

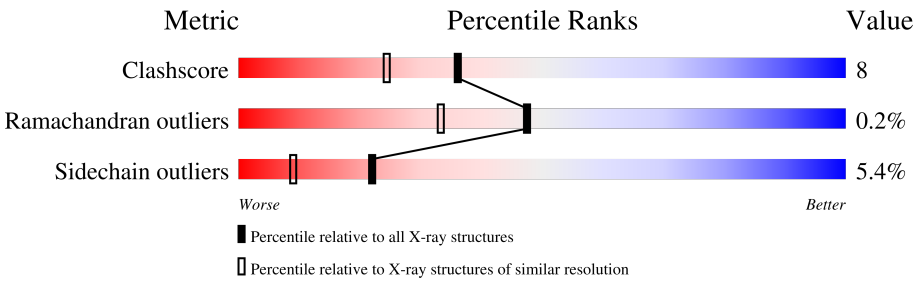
MolProbity	:	4.02b-467
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

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 1.80 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	141614	6793 (1.80-1.80)
Ramachandran outliers	138981	6697 (1.80-1.80)
Sidechain outliers	138945	6696 (1.80-1.80)

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%

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	571	83% 14% • •

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
4	FUM	A	806	-	X	-	-

2 Entry composition (i)

There are 7 unique types of molecules in this entry. The entry contains 5057 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.

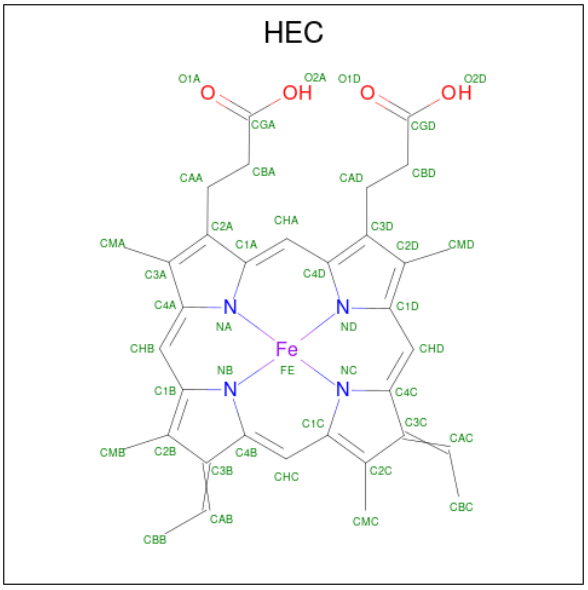
- Molecule 1 is a protein called FUMARATE REDUCTASE FLAVOPROTEIN SUBUNIT.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	568	4212	2612	742	833	25	0	0	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	365	ALA	HIS	engineered mutation	UNP Q07WU7

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



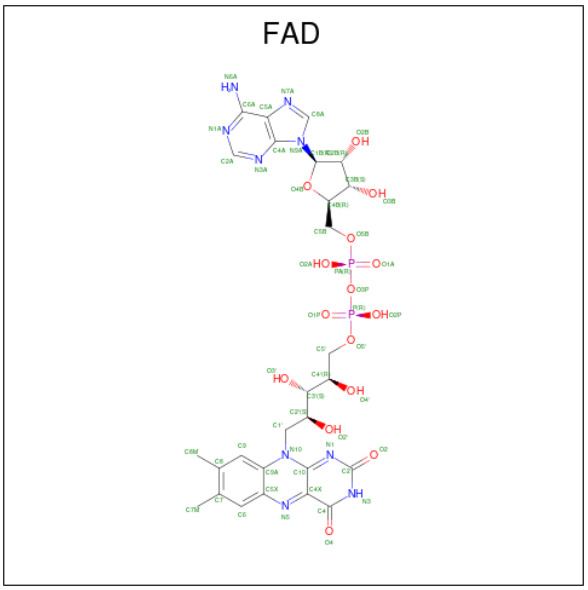
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	Fe	N	O		
2	A	1	43	34	1	4	4	0	0
			Total	C	Fe	N	O		
2	A	1	43	34	1	4	4	0	0
			Total	C	Fe	N	O		
2	A	1	43	34	1	4	4	0	0

Continued on next page...

Continued from previous page...

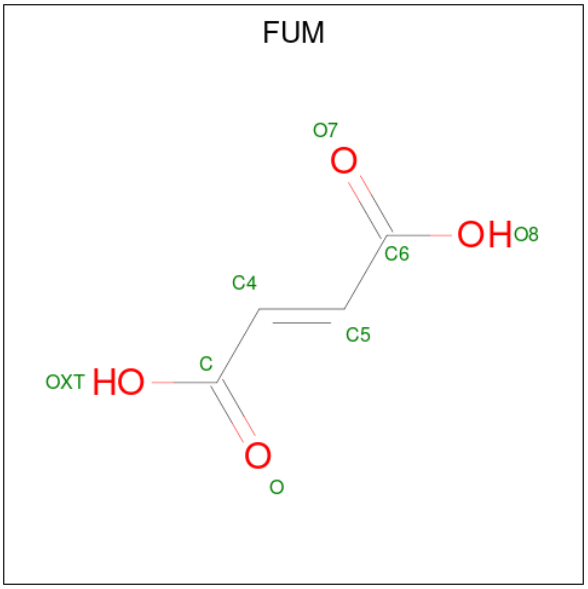
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	Fe	N	O		
2	A	1	43	34	1	4	4	0	0

- Molecule 3 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula: $C_{27}H_{33}N_9O_{15}P_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
3	A	1	53	27	9	15	2	0	0

- Molecule 4 is FUMARIC ACID (three-letter code: FUM) (formula: $C_4H_4O_4$).

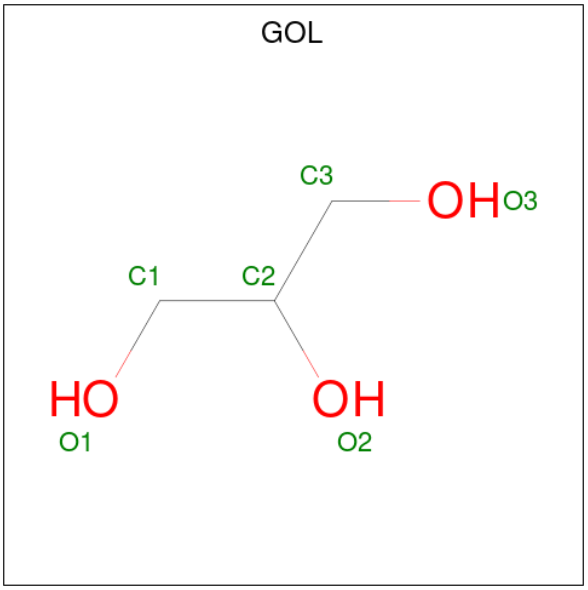


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
4	A	1	8	4	4	0	0

- Molecule 5 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Na		
5	A	1	1	1	0	0

- Molecule 6 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



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

- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
7	A	605	605	605	0	0

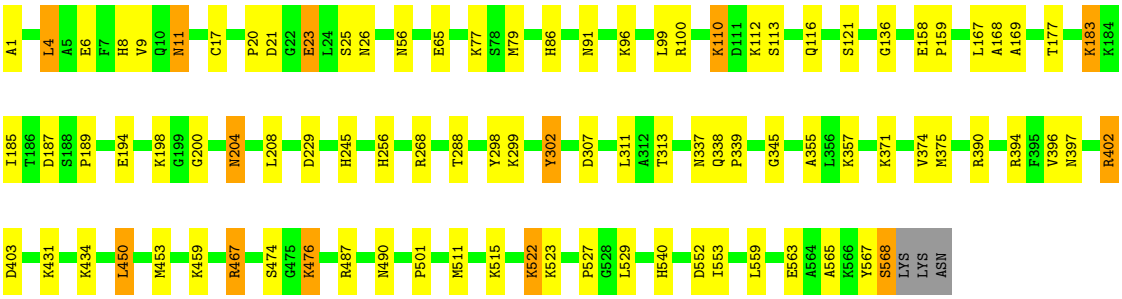
3 Residue-property plots (i)

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. 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. 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.

Note EDS was not executed.

• Molecule 1: FUMARATE REDUCTASE FLAVOPROTEIN SUBUNIT

Chain A: 83% 14% ••



4 Data and refinement statistics (i)

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	45.57Å 92.17Å 78.49Å 90.00° 91.08° 90.00°	Depositor
Resolution (Å)	18.00 – 1.80	Depositor
% Data completeness (in resolution range)	86.0 (18.00-1.80)	Depositor
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	SHELXL-97	Depositor
R, R_{free}	0.181 , 0.247	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	5057	wwPDB-VP
Average B, all atoms (Å^2)	21.0	wwPDB-VP

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: GOL, HEC, FAD, FUM, NA

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	0.36	0/4281	1.00	12/5787 (0.2%)

There are no bond length outliers.

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	187	ASP	CB-CG-OD1	7.75	125.28	118.30
1	A	86	HIS	CA-CB-CG	7.69	126.67	113.60
1	A	402	ARG	NE-CZ-NH2	-6.12	117.24	120.30
1	A	394	ARG	NE-CZ-NH2	-5.97	117.31	120.30
1	A	100	ARG	NE-CZ-NH1	5.85	123.22	120.30
1	A	567	TYR	CB-CG-CD1	-5.61	117.64	121.00
1	A	302	TYR	CB-CG-CD1	5.25	124.15	121.00
1	A	298	TYR	CB-CG-CD2	5.23	124.14	121.00
1	A	390	ARG	NE-CZ-NH2	-5.21	117.70	120.30
1	A	402	ARG	NE-CZ-NH1	5.09	122.85	120.30
1	A	268	ARG	NE-CZ-NH1	5.04	122.82	120.30
1	A	567	TYR	C-N-CA	5.03	134.28	121.70

There are no chirality outliers.

There are no planarity outliers.

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	4212	0	4150	67	0
2	A	172	0	120	6	0
3	A	53	0	31	1	0
4	A	8	0	1	2	0
5	A	1	0	0	0	0
6	A	6	0	8	1	0
7	A	605	0	0	18	0
All	All	5057	0	4310	70	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 8.

All (70) 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:431:LYS:HD3	1:A:434:LYS:HD3	1.48	0.95
1:A:204:ASN:HD22	1:A:204:ASN:H	1.29	0.79
1:A:200:GLY:HA3	1:A:204:ASN:HD21	1.47	0.77
1:A:229:ASP:H	1:A:256:HIS:HE1	1.33	0.77
1:A:112:LYS:HD2	7:A:2207:HOH:O	1.84	0.75
1:A:9:VAL:HG13	7:A:2020:HOH:O	1.89	0.73
1:A:450:LEU:HD23	1:A:453:MET:HE2	1.71	0.71
1:A:65:GLU:HG2	7:A:2108:HOH:O	1.92	0.67
1:A:431:LYS:HA	1:A:434:LYS:HD2	1.77	0.67
1:A:375:MET:HE3	4:A:806:FUM:O8	1.98	0.64
1:A:1:ALA:HA	1:A:6:GLU:OE2	1.99	0.63
1:A:21:ASP:OD1	1:A:23:GLU:HB3	1.99	0.63
1:A:229:ASP:H	1:A:256:HIS:CE1	2.15	0.62
1:A:189:PRO:HG2	7:A:2284:HOH:O	1.99	0.62
1:A:396:VAL:HG22	7:A:2465:HOH:O	2.01	0.61
1:A:371:LYS:HD2	6:A:850:GOL:H11	1.83	0.60
1:A:11:ASN:HB3	7:A:2017:HOH:O	2.01	0.60
2:A:801:HEC:HMC1	2:A:801:HEC:HBC3	1.84	0.59
1:A:17:CYS:HB3	7:A:2025:HOH:O	2.02	0.58
1:A:20:PRO:HD3	7:A:2052:HOH:O	2.04	0.57
1:A:565:ALA:O	1:A:568:SER:HA	2.06	0.56
1:A:307:ASP:HA	7:A:2390:HOH:O	2.06	0.56
2:A:802:HEC:HMC1	2:A:802:HEC:HBC3	1.87	0.55
1:A:515:LYS:HE3	1:A:563:GLU:OE1	2.06	0.55
1:A:112:LYS:O	1:A:116:GLN:HG3	2.08	0.54
1:A:450:LEU:HD23	1:A:453:MET:CE	2.38	0.53
1:A:4:LEU:HD22	1:A:8:HIS:CE1	2.44	0.53

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:476:LYS:HG2	1:A:476:LYS:O	2.08	0.52
1:A:23:GLU:HG2	1:A:23:GLU:O	2.06	0.51
1:A:110:LYS:N	1:A:110:LYS:HD2	2.22	0.51
1:A:374:VAL:HB	2:A:804:HEC:HMD3	1.93	0.51
1:A:540:HIS:HD2	7:A:2573:HOH:O	1.93	0.51
1:A:77:LYS:HG2	7:A:2128:HOH:O	2.10	0.51
1:A:501:PRO:HD2	7:A:2548:HOH:O	2.11	0.51
1:A:177:THR:OG1	1:A:245:HIS:HE1	1.93	0.50
1:A:136:GLY:HA3	1:A:553:ILE:HD12	1.95	0.49
2:A:803:HEC:HBC3	2:A:803:HEC:HMC1	1.95	0.49
1:A:20:PRO:HD2	7:A:2032:HOH:O	2.13	0.48
1:A:183:LYS:O	1:A:185:ILE:HD12	2.13	0.48
1:A:302:TYR:HA	7:A:2384:HOH:O	2.13	0.48
1:A:431:LYS:HD3	1:A:434:LYS:CD	2.34	0.48
1:A:523:LYS:HE3	7:A:2562:HOH:O	2.14	0.47
1:A:204:ASN:H	1:A:204:ASN:ND2	2.03	0.47
1:A:168:ALA:HA	3:A:805:FAD:N5	2.29	0.47
1:A:540:HIS:HE1	1:A:552:ASP:OD2	1.98	0.46
1:A:4:LEU:HD22	1:A:8:HIS:HE1	1.81	0.46
1:A:397:ASN:HB3	7:A:2465:HOH:O	2.16	0.46
1:A:357:LYS:HE3	1:A:357:LYS:HB2	1.78	0.45
1:A:169:ALA:HB1	1:A:431:LYS:HE3	1.97	0.45
1:A:522:LYS:O	1:A:523:LYS:HB2	2.16	0.45
1:A:77:LYS:NZ	7:A:2129:HOH:O	2.50	0.45
1:A:355:ALA:HB3	1:A:511:MET:HE3	1.98	0.44
1:A:403:ASP:N	1:A:403:ASP:OD1	2.50	0.44
1:A:553:ILE:HD13	1:A:553:ILE:HG21	1.80	0.43
1:A:169:ALA:CB	1:A:431:LYS:HE3	2.48	0.43
1:A:79:MET:CE	1:A:96:LYS:HE3	2.49	0.43
1:A:229:ASP:N	1:A:256:HIS:HE1	2.09	0.42
1:A:402:ARG:HH22	4:A:806:FUM:C6	2.31	0.42
1:A:453:MET:HE2	1:A:453:MET:HB2	1.80	0.42
1:A:467:ARG:NE	7:A:2522:HOH:O	2.52	0.42
1:A:158:GLU:HB3	1:A:159:PRO:HD2	2.01	0.42
1:A:313:THR:OG1	1:A:345:GLY:HA3	2.20	0.42
1:A:338:GLN:HB2	1:A:339:PRO:HD2	2.01	0.42
1:A:167:LEU:HD11	2:A:804:HEC:HBC2	2.02	0.42
1:A:474:SER:O	1:A:476:LYS:HD3	2.20	0.41
1:A:288:THR:HB	1:A:527:PRO:HB2	2.02	0.41
1:A:311:LEU:HD21	1:A:529:LEU:HD11	2.01	0.41
1:A:474:SER:OG	1:A:476:LYS:NZ	2.52	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:25:SER:O	1:A:26:ASN:HB3	2.20	0.41
1:A:167:LEU:CD1	2:A:804:HEC:HBC2	2.51	0.40

There are no symmetry-related clashes.

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	566/571 (99%)	551 (97%)	14 (2%)	1 (0%)	47 33

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	113	SER

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	441/444 (99%)	417 (95%)	24 (5%)	22 9

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	LEU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	11	ASN
1	A	23	GLU
1	A	56	ASN
1	A	91	ASN
1	A	99	LEU
1	A	110	LYS
1	A	121	SER
1	A	183	LYS
1	A	194	GLU
1	A	198	LYS
1	A	204	ASN
1	A	208	LEU
1	A	299	LYS
1	A	337	ASN
1	A	450	LEU
1	A	459	LYS
1	A	467	ARG
1	A	476	LYS
1	A	487	ARG
1	A	490	ASN
1	A	522	LYS
1	A	559	LEU
1	A	568	SER

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	56	ASN
1	A	91	ASN
1	A	204	ASN
1	A	245	HIS
1	A	256	HIS
1	A	490	ASN
1	A	540	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 monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 8 ligands modelled in this entry, 1 is monoatomic - leaving 7 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	$\#\|Z\| > 2$
3	FAD	A	805	-	53,58,58	1.34	6 (11%)	68,89,89	1.31	7 (10%)
2	HEC	A	803	1	32,50,50	1.61	2 (6%)	24,82,82	1.41	2 (8%)
2	HEC	A	802	1	32,50,50	1.45	2 (6%)	24,82,82	1.23	2 (8%)
2	HEC	A	804	1	32,50,50	1.59	2 (6%)	24,82,82	1.36	2 (8%)
4	FUM	A	806	-	7,7,7	2.74	4 (57%)	8,8,8	1.53	1 (12%)
6	GOL	A	850	-	5,5,5	0.35	0	5,5,5	0.57	0
2	HEC	A	801	1	32,50,50	1.46	2 (6%)	24,82,82	1.44	3 (12%)

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
3	FAD	A	805	-	-	7/30/50/50	0/6/6/6
2	HEC	A	803	1	-	5/10/54/54	-
2	HEC	A	802	1	-	2/10/54/54	-
2	HEC	A	804	1	-	5/10/54/54	-
4	FUM	A	806	-	-	3/5/5/5	-

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	GOL	A	850	-	-	3/4/4/4	-
2	HEC	A	801	1	-	3/10/54/54	-

All (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	804	HEC	C2B-C3B	-5.18	1.35	1.40
2	A	803	HEC	C2B-C3B	-4.90	1.35	1.40
2	A	803	HEC	C3C-C2C	-4.80	1.35	1.40
2	A	804	HEC	C3C-C2C	-4.78	1.35	1.40
2	A	801	HEC	C2B-C3B	-4.70	1.35	1.40
2	A	802	HEC	C3C-C2C	-4.48	1.36	1.40
2	A	802	HEC	C2B-C3B	-4.35	1.36	1.40
4	A	806	FUM	O-C	4.30	1.34	1.23
2	A	801	HEC	C3C-C2C	-4.02	1.36	1.40
4	A	806	FUM	O7-C6	3.67	1.32	1.23
3	A	805	FAD	C9-C8	3.61	1.44	1.39
3	A	805	FAD	C2'-C3'	3.45	1.60	1.53
3	A	805	FAD	C6-C5X	3.36	1.45	1.40
4	A	806	FUM	OXT-C	-3.34	1.21	1.30
4	A	806	FUM	O8-C6	-2.97	1.22	1.30
3	A	805	FAD	C1'-C2'	-2.88	1.48	1.52
3	A	805	FAD	O4-C4	-2.76	1.18	1.23
3	A	805	FAD	C5'-C4'	2.09	1.54	1.51

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	A	805	FAD	O5'-C5'-C4'	-5.94	93.49	109.36
3	A	805	FAD	O4B-C1B-C2B	-4.29	100.66	106.93
2	A	802	HEC	CBA-CAA-C2A	3.75	118.92	112.60
2	A	803	HEC	CMC-C2C-C1C	-3.70	122.78	128.46
2	A	803	HEC	CMC-C2C-C3C	3.69	130.16	125.82
2	A	801	HEC	CMB-C2B-C1B	-3.62	122.91	128.46
2	A	801	HEC	CMB-C2B-C3B	3.19	129.57	125.82
2	A	801	HEC	CMD-C2D-C1D	-3.18	123.57	128.46
2	A	804	HEC	CBA-CAA-C2A	3.01	117.67	112.60
2	A	804	HEC	CMC-C2C-C1C	-2.72	124.28	128.46
4	A	806	FUM	C5-C4-C	-2.67	113.55	127.26
3	A	805	FAD	O4'-C4'-C3'	-2.61	102.75	109.10
3	A	805	FAD	O2'-C2'-C3'	-2.32	103.47	109.10
3	A	805	FAD	C6-C5X-C9A	2.25	122.13	118.94

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	A	805	FAD	O2P-P-O1P	2.18	123.02	112.24
3	A	805	FAD	O2A-PA-O1A	2.03	122.26	112.24
2	A	802	HEC	CMB-C2B-C1B	-2.02	125.36	128.46

There are no chirality outliers.

All (28) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	801	HEC	C4D-C3D-CAD-CBD
2	A	803	HEC	C3A-C2A-CAA-CBA
3	A	805	FAD	C2'-C3'-C4'-O4'
3	A	805	FAD	C2'-C3'-C4'-C5'
3	A	805	FAD	O3'-C3'-C4'-C5'
4	A	806	FUM	C-C4-C5-C6
3	A	805	FAD	O3'-C3'-C4'-O4'
4	A	806	FUM	C4-C5-C6-O8
4	A	806	FUM	C4-C5-C6-O7
6	A	850	GOL	O1-C1-C2-O2
6	A	850	GOL	O1-C1-C2-C3
2	A	803	HEC	C3D-CAD-CBD-CGD
6	A	850	GOL	O2-C2-C3-O3
2	A	802	HEC	CAD-CBD-CGD-O2D
2	A	803	HEC	C2A-CAA-CBA-CGA
2	A	804	HEC	C2A-CAA-CBA-CGA
2	A	803	HEC	CAA-CBA-CGA-O2A
2	A	804	HEC	CAA-CBA-CGA-O2A
2	A	804	HEC	CAA-CBA-CGA-O1A
2	A	803	HEC	CAA-CBA-CGA-O1A
2	A	802	HEC	CAD-CBD-CGD-O1D
3	A	805	FAD	P-O3P-PA-O1A
2	A	801	HEC	CAD-CBD-CGD-O1D
2	A	804	HEC	CAD-CBD-CGD-O1D
2	A	801	HEC	CAD-CBD-CGD-O2D
3	A	805	FAD	P-O3P-PA-O2A
3	A	805	FAD	O4B-C4B-C5B-O5B
2	A	804	HEC	CAD-CBD-CGD-O2D

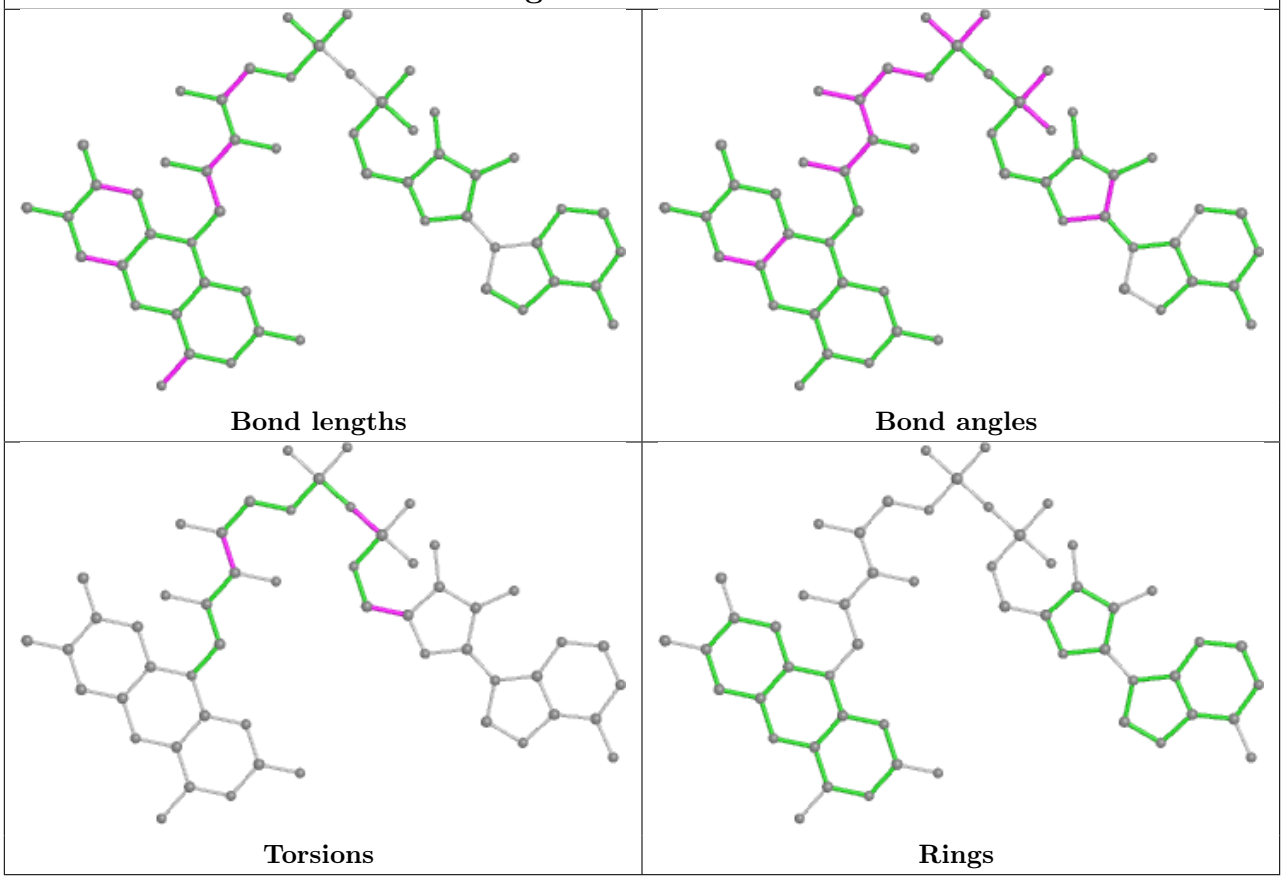
There are no ring outliers.

7 monomers are involved in 10 short contacts:

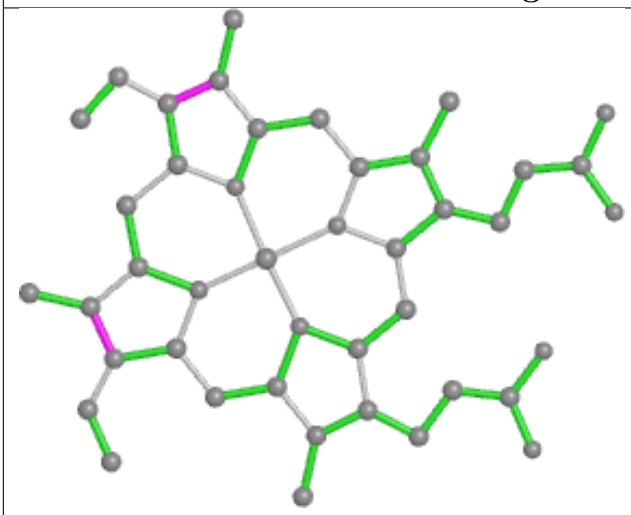
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	805	FAD	1	0
2	A	803	HEC	1	0
2	A	802	HEC	1	0
2	A	804	HEC	3	0
4	A	806	FUM	2	0
6	A	850	GOL	1	0
2	A	801	HEC	1	0

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.

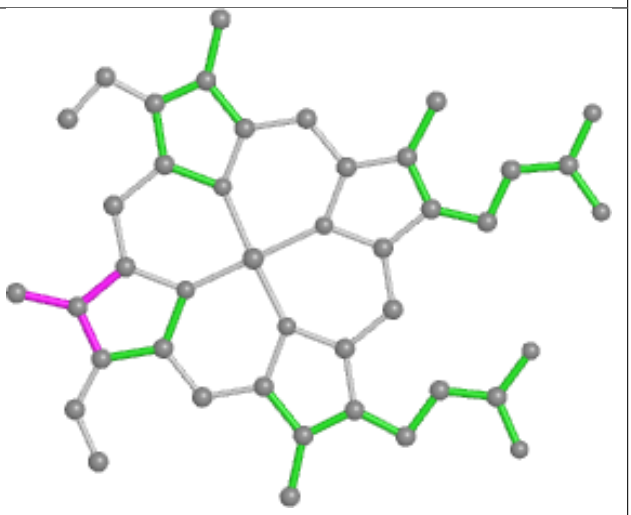
Ligand FAD A 805



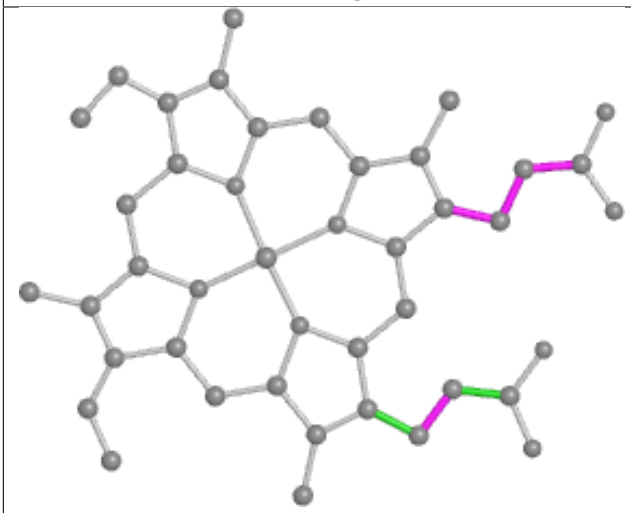
Ligand HEC A 803



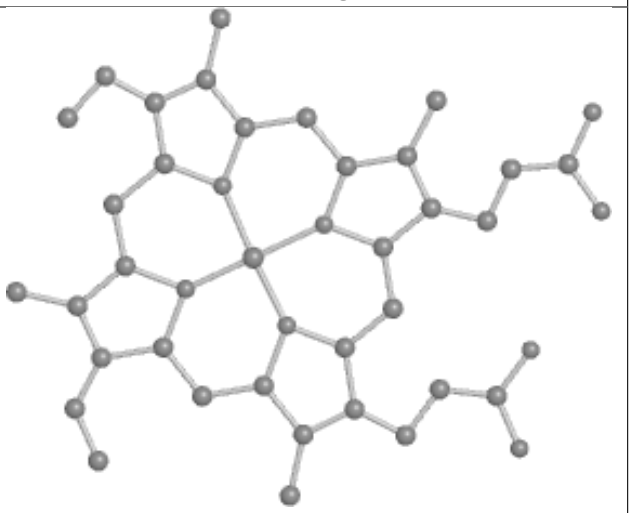
Bond lengths



Bond angles

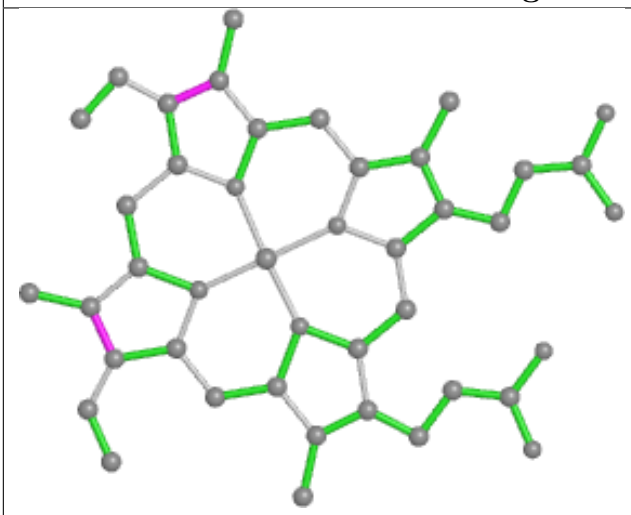


Torsions

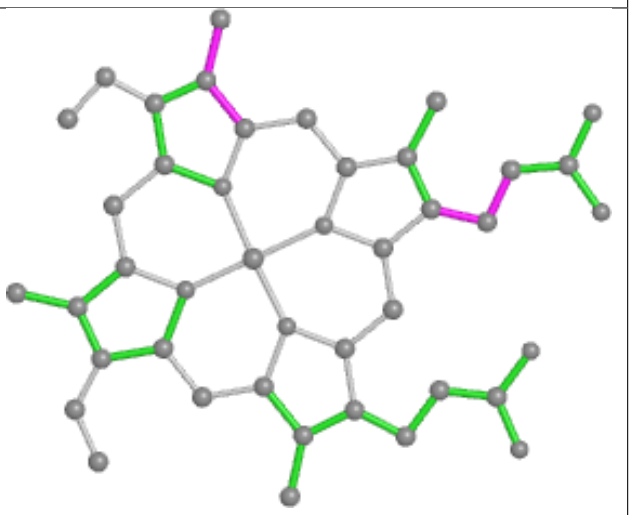


Rings

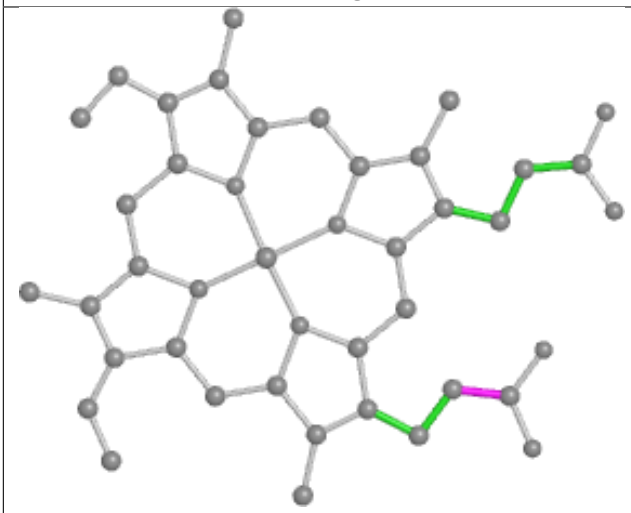
Ligand HEC A 802



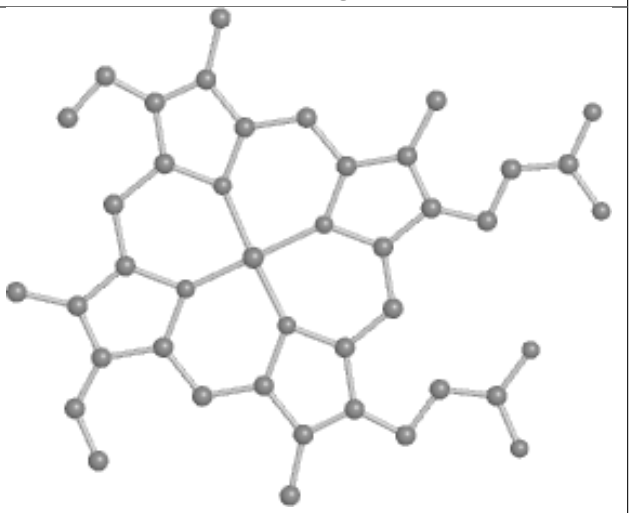
Bond lengths



Bond angles

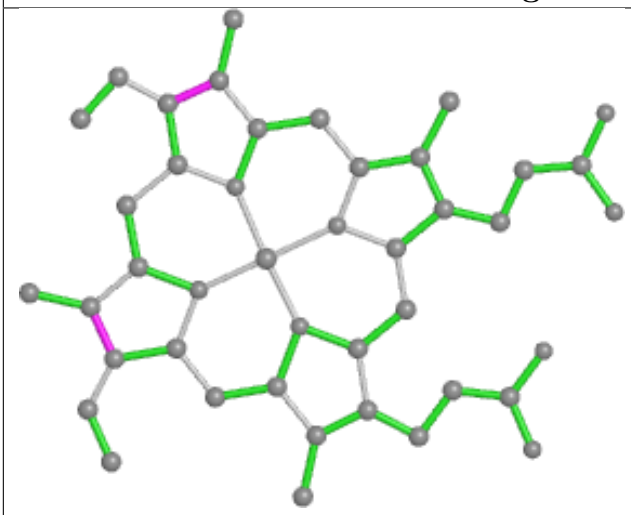


Torsions

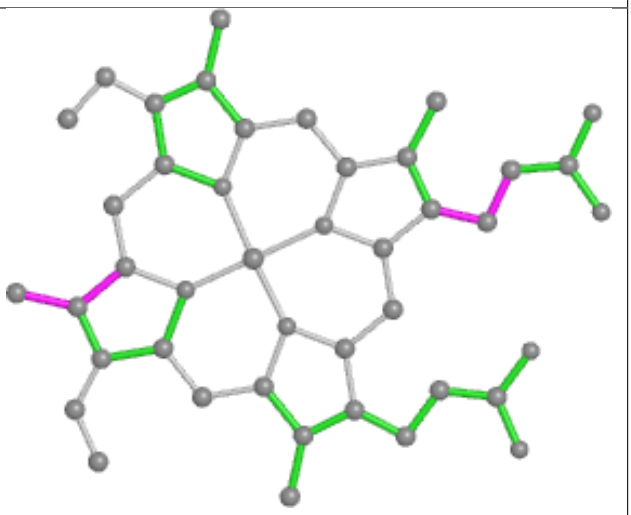


Rings

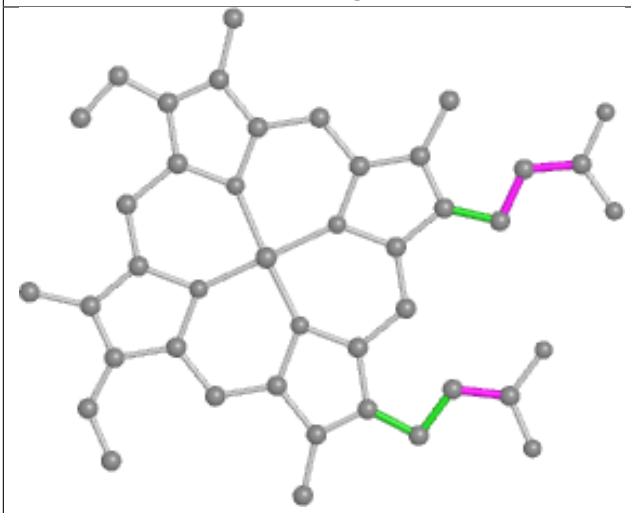
Ligand HEC A 804



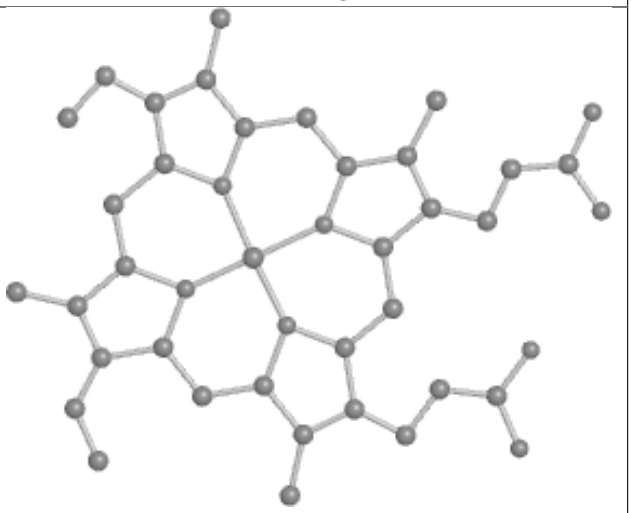
Bond lengths



Bond angles

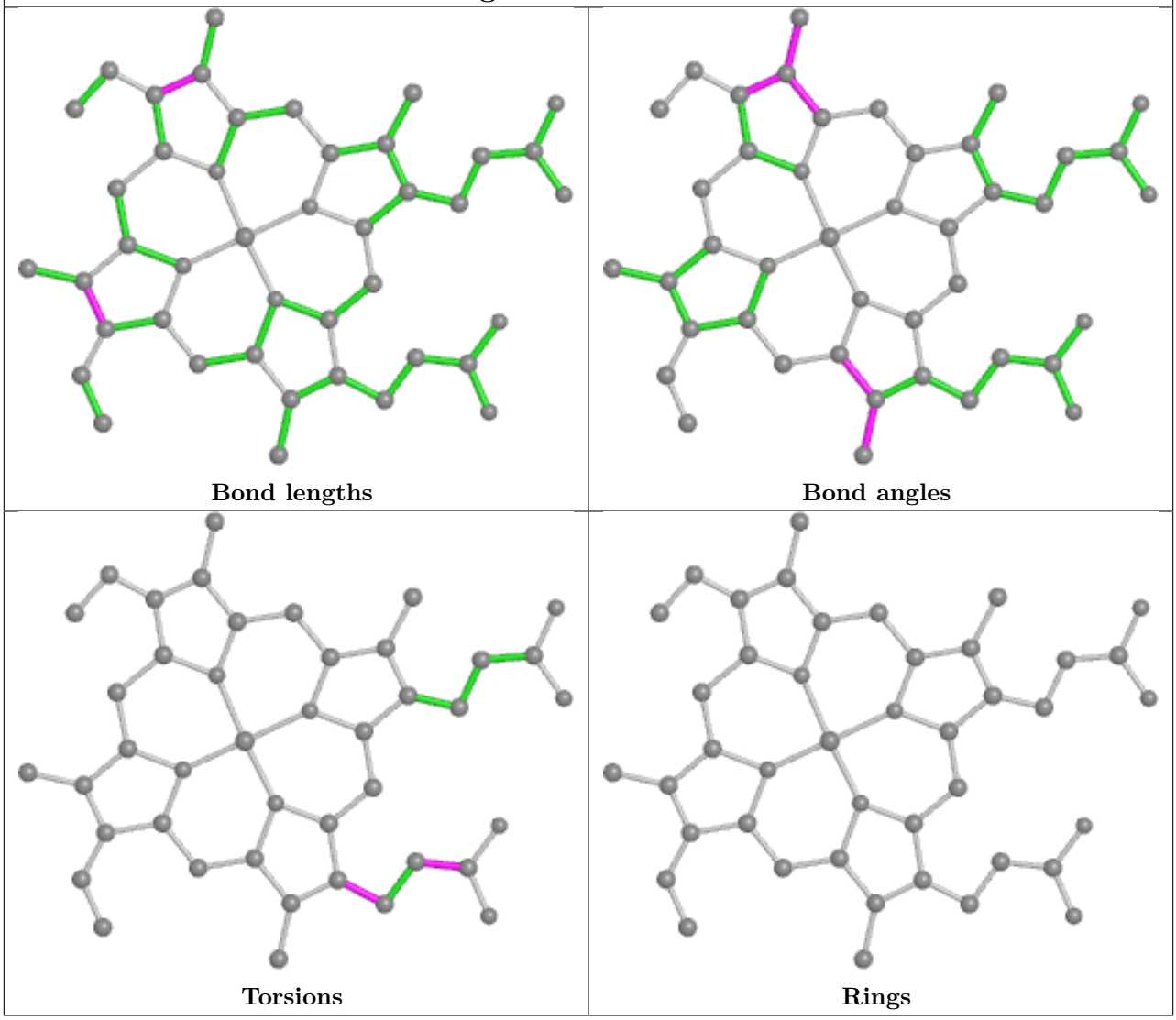


Torsions



Rings

Ligand HEC A 801



5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data (i)

6.1 Protein, DNA and RNA chains (i)

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates (i)

EDS was not executed - this section is therefore empty.

6.4 Ligands (i)

EDS was not executed - this section is therefore empty.

6.5 Other polymers (i)

EDS was not executed - this section is therefore empty.