



Full wwPDB X-ray Structure Validation Report ⓘ

May 29, 2020 – 01:09 am BST

PDB ID : 1AOF
Title : CYTOCHROME CD1 NITRITE REDUCTASE, REDUCED FORM
Authors : Williams, P.A.; Fulop, V.
Deposited on : 1997-07-02
Resolution : 2.00 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

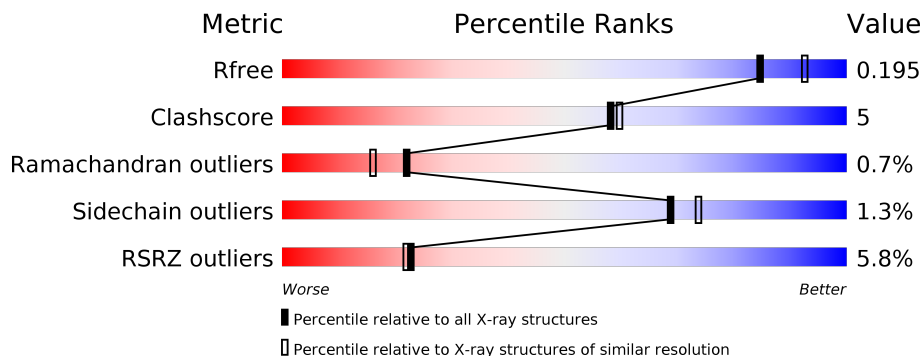
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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}	130704	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

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 on 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	567	 8% 83% 10% • 6%
1	B	567	 3% 84% 11% •

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 9085 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.

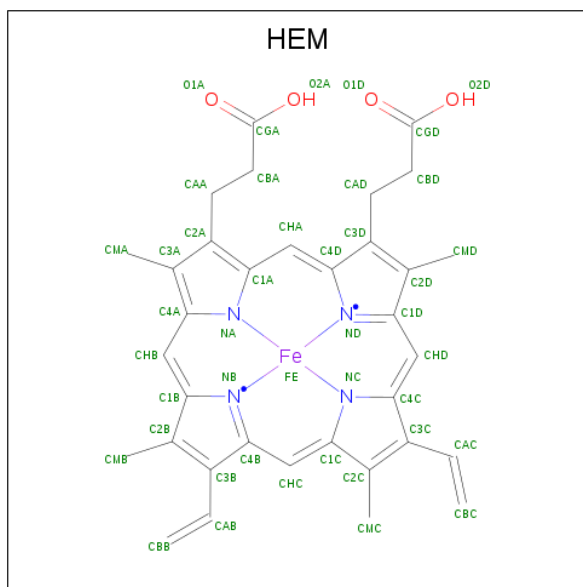
- Molecule 1 is a protein called NITRITE REDUCTASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	532	4133	2618	689	811	15	0	0	0
1	B	542	4210	2664	702	829	15	0	0	0

There are 8 discrepancies between the modelled and reference sequences:

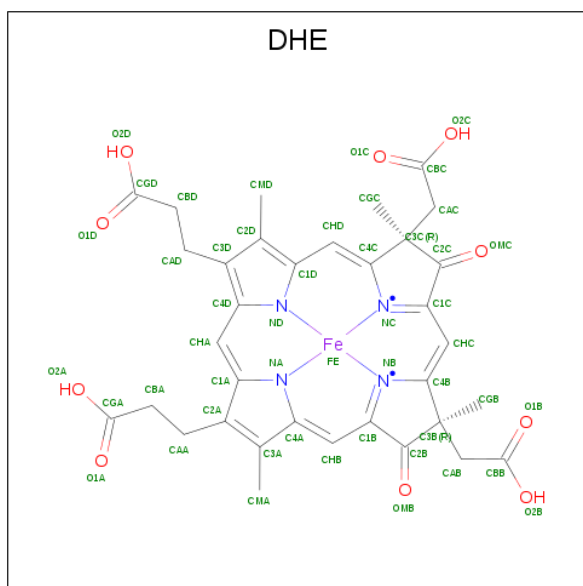
Chain	Residue	Modelled	Actual	Comment	Reference
A	160	MET	GLU	CONFLICT	UNP P72181
A	185	SER	THR	CONFLICT	UNP P72181
A	191	THR	SER	CONFLICT	UNP P72181
A	331	ASN	ASP	CONFLICT	UNP P72181
B	160	MET	GLU	CONFLICT	UNP P72181
B	185	SER	THR	CONFLICT	UNP P72181
B	191	THR	SER	CONFLICT	UNP P72181
B	331	ASN	ASP	CONFLICT	UNP P72181

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



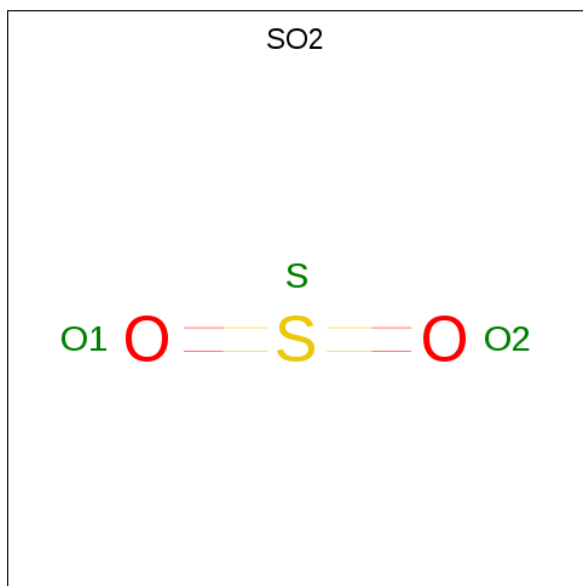
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
2	A	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
2	B	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		

- Molecule 3 is HEME D (three-letter code: DHE) (formula: $C_{34}H_{32}FeN_4O_{10}$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
3	A	1	Total	C	Fe	N	O	0	0
			49	34	1	4	10		
3	B	1	Total	C	Fe	N	O	0	0
			49	34	1	4	10		

- Molecule 4 is SULFUR DIOXIDE (three-letter code: SO2) (formula: O₂S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	S	0	0
			3	2	1		
4	B	1	Total	O	S	0	0
			3	2	1		

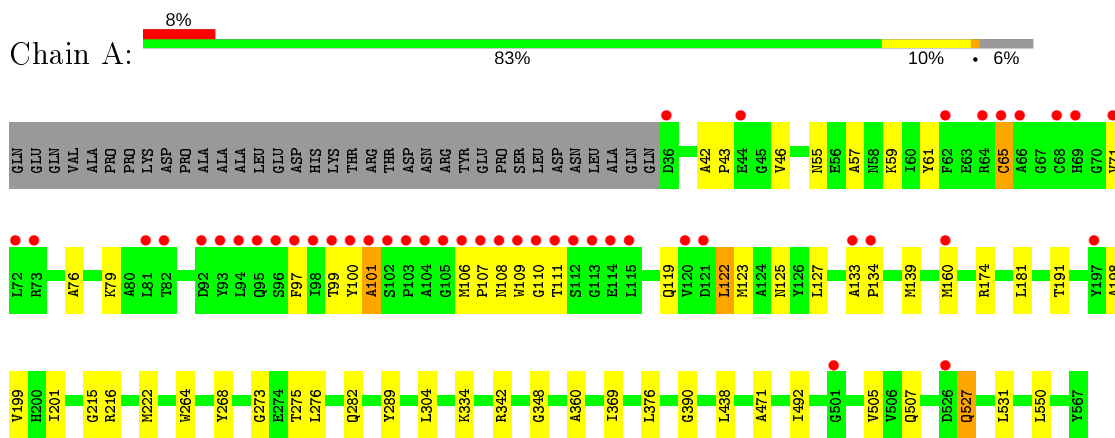
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	267	Total	O	0	0
			267	267		
5	B	285	Total	O	0	0
			285	285		

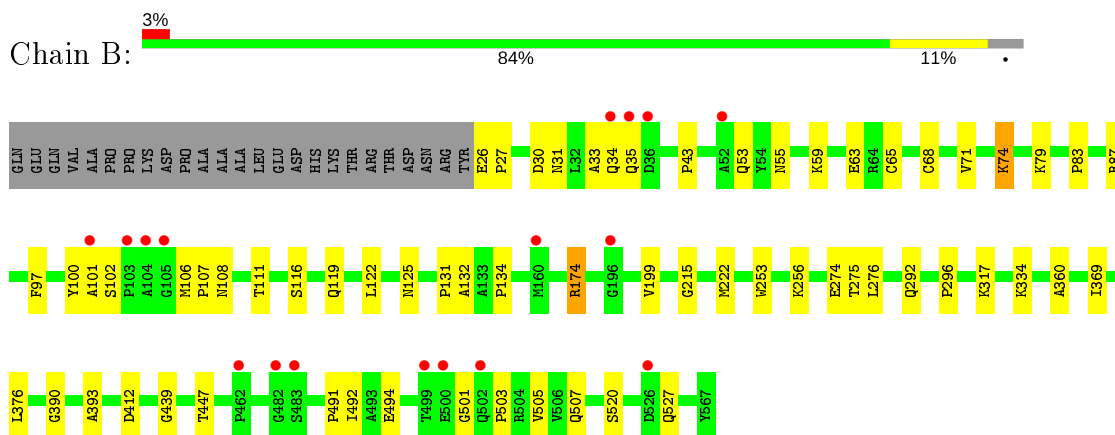
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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: NITRITE REDUCTASE



- Molecule 1: NITRITE REDUCTASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	107.40Å 61.70Å 101.20Å 90.00° 112.10° 90.00°	Depositor
Resolution (Å)	20.00 – 2.00 19.90 – 2.00	Depositor EDS
% Data completeness (in resolution range)	94.0 (20.00-2.00) 97.1 (19.90-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.53 (at 2.01Å)	Xtrriage
Refinement program	X-PLOR 3.8	Depositor
R, R_{free}	0.170 , 0.201 0.162 , 0.195	Depositor DCC
R_{free} test set	3245 reflections (4.02%)	wwPDB-VP
Wilson B-factor (Å ²)	21.6	Xtrriage
Anisotropy	0.260	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 55.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	9085	wwPDB-VP
Average B, all atoms (Å ²)	26.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.17% of the height of the origin peak. No significant pseudotranslation is detected.*

¹ Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: HEM, DHE, SO2

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	0.38	0/4235	0.68	2/5762 (0.0%)
1	B	0.36	0/4313	0.68	2/5869 (0.0%)
All	All	0.37	0/8548	0.68	4/11631 (0.0%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	390	GLY	N-CA-C	-6.77	96.18	113.10
1	B	390	GLY	N-CA-C	-6.71	96.32	113.10
1	A	360	ALA	N-CA-C	-5.93	95.00	111.00
1	B	360	ALA	N-CA-C	-5.64	95.78	111.00

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	4133	0	3992	41	0
1	B	4210	0	4063	41	0
2	A	43	0	30	2	0
2	B	43	0	30	1	0
3	A	49	0	24	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	49	0	24	1	0
4	A	3	0	0	0	0
4	B	3	0	0	0	0
5	A	267	0	0	0	0
5	B	285	0	0	1	0
All	All	9085	0	8163	80	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 5.

All (80) 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:57:ALA:HB2	1:A:122:LEU:HD22	1.59	0.84
1:A:334:LYS:HE2	1:B:334:LYS:HE2	1.57	0.84
1:B:101:ALA:HB1	1:B:106:MET:O	1.82	0.79
1:A:101:ALA:HB1	1:A:106:MET:O	1.87	0.75
1:B:101:ALA:HB2	1:B:108:ASN:N	2.02	0.74
1:A:99:THR:HG23	1:A:110:GLY:HA3	1.68	0.74
1:B:101:ALA:HB2	1:B:108:ASN:CA	2.18	0.73
1:B:26:GLU:HG3	1:B:68:CYS:SG	2.29	0.72
1:B:199:VAL:HA	1:B:215:GLY:HA2	1.77	0.64
1:B:116:SER:H	1:B:119:GLN:HE21	1.47	0.61
1:A:100:TYR:O	1:A:101:ALA:HB3	1.99	0.61
1:A:101:ALA:HB2	1:A:107:PRO:C	2.22	0.60
1:B:100:TYR:O	1:B:101:ALA:HB3	2.00	0.60
1:B:501:GLY:O	1:B:503:PRO:HD3	2.02	0.59
1:B:101:ALA:HB2	1:B:108:ASN:HA	1.84	0.58
1:A:101:ALA:HB2	1:A:108:ASN:CA	2.35	0.56
1:A:71:VAL:HG12	1:A:134:PRO:CG	2.36	0.56
1:A:101:ALA:HB2	1:A:108:ASN:N	2.21	0.56
1:B:71:VAL:CG1	1:B:134:PRO:HG2	2.36	0.56
1:B:101:ALA:HB2	1:B:107:PRO:C	2.25	0.55
1:B:74:LYS:HG2	1:B:296:PRO:HG3	1.88	0.54
1:A:101:ALA:HB2	1:A:108:ASN:HA	1.89	0.54
1:A:119:GLN:O	1:A:123:MET:HG2	2.07	0.54
1:A:43:PRO:HD2	1:A:46:VAL:HB	1.90	0.53
1:B:79:LYS:HG3	2:B:601:HEM:HAD2	1.91	0.53
1:B:53:GLN:HB3	1:B:122:LEU:HD21	1.91	0.52
1:B:31:ASN:O	1:B:34:GLN:HG2	2.10	0.52
1:B:222:MET:HE1	1:B:274:GLU:HA	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:253:TRP:CE3	1:B:256:LYS:HD2	2.46	0.51
1:B:55:ASN:O	1:B:59:LYS:HG3	2.11	0.51
1:B:27:PRO:HA	3:B:602:DHE:HAB2	1.92	0.50
1:A:61:TYR:CE1	1:A:65:CYS:HB2	2.46	0.50
1:A:222:MET:HE1	1:A:273:GLY:O	2.12	0.49
1:A:527:GLN:HE21	1:A:527:GLN:HA	1.77	0.49
1:B:116:SER:H	1:B:119:GLN:NE2	2.10	0.49
1:A:108:ASN:HB2	1:A:111:THR:OG1	2.12	0.48
1:B:174:ARG:HG2	1:B:174:ARG:HH11	1.77	0.48
1:A:304:LEU:HD21	1:A:348:GLY:HA2	1.95	0.48
1:A:61:TYR:HA	1:A:65:CYS:SG	2.53	0.48
1:B:35:GLN:HB3	5:B:721:HOH:O	2.12	0.48
1:B:43:PRO:HG3	1:B:87:ARG:CZ	2.44	0.48
1:A:369:ILE:HD13	1:A:376:LEU:HA	1.95	0.48
1:B:132:ALA:O	1:B:134:PRO:HD3	2.14	0.47
1:A:275:THR:O	1:A:276:LEU:HB2	2.16	0.46
1:A:123:MET:O	1:A:127:LEU:HG	2.16	0.46
1:B:43:PRO:HG3	1:B:87:ARG:NH2	2.31	0.46
1:A:198:ALA:HB3	1:A:216:ARG:HD3	1.99	0.45
1:A:42:ALA:HA	1:A:43:PRO:HD3	1.81	0.45
1:A:438:LEU:HG	1:A:471:ALA:HB2	1.99	0.45
1:A:492:ILE:HB	1:A:505:VAL:HG21	1.99	0.45
1:B:33:ALA:CB	1:B:63:GLU:HG3	2.47	0.45
1:B:83:PRO:HG2	1:B:131:PRO:HD3	1.98	0.45
1:A:100:TYR:O	1:A:101:ALA:CB	2.63	0.45
1:A:71:VAL:HG12	1:A:134:PRO:HG3	1.98	0.45
1:A:79:LYS:HG2	2:A:601:HEM:HAD2	1.99	0.44
1:A:199:VAL:HA	1:A:215:GLY:HA2	1.99	0.44
1:A:134:PRO:HB2	1:A:264:TRP:CZ2	2.52	0.44
1:B:33:ALA:O	1:B:59:LYS:HG2	2.18	0.43
1:B:102:SER:HB3	1:B:106:MET:HB3	2.00	0.43
1:A:201:ILE:HD13	3:A:602:DHE:O2A	2.18	0.43
1:B:393:ALA:HB1	1:B:447:THR:HG22	2.00	0.43
1:A:289:TYR:HE1	1:A:342:ARG:HG3	1.83	0.43
1:B:275:THR:O	1:B:276:LEU:HB2	2.19	0.43
1:A:531:LEU:HG	1:A:550:LEU:HD21	1.99	0.43
1:B:492:ILE:HB	1:B:505:VAL:HG21	2.00	0.43
1:A:106:MET:HE1	1:A:109:TRP:CD1	2.54	0.43
1:B:369:ILE:HD13	1:B:376:LEU:HA	2.01	0.42
1:B:491:PRO:HB2	1:B:494:GLU:HB3	2.00	0.42
1:B:505:VAL:HA	1:B:520:SER:O	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:133:ALA:HA	1:A:134:PRO:HD3	1.89	0.42
1:B:412:ASP:HB3	1:B:439:GLY:HA2	2.00	0.41
1:A:268:TYR:CZ	1:A:282:GLN:HB2	2.56	0.41
1:B:108:ASN:HB2	1:B:111:THR:OG1	2.20	0.41
1:A:106:MET:HB2	2:A:601:HEM:C4D	2.56	0.41
1:A:139:MET:HB2	1:B:292:GLN:HB2	2.01	0.41
1:A:181:LEU:HD12	1:A:191:THR:CG2	2.51	0.41
1:B:106:MET:HA	1:B:107:PRO:HD3	1.88	0.41
1:A:99:THR:HA	1:A:110:GLY:H	1.85	0.41
1:A:55:ASN:O	1:A:59:LYS:HG3	2.21	0.41
1:B:174:ARG:HG2	1:B:174:ARG:NH1	2.36	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	530/567 (94%)	505 (95%)	21 (4%)	4 (1%)	19 13
1	B	540/567 (95%)	517 (96%)	20 (4%)	3 (1%)	25 19
All	All	1070/1134 (94%)	1022 (96%)	41 (4%)	7 (1%)	22 16

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	101	ALA
1	A	507	GLN
1	B	507	GLN
1	A	174	ARG
1	B	174	ARG
1	A	76	ALA
1	B	317	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	440/470 (94%)	434 (99%)	6 (1%)	67	72
1	B	449/470 (96%)	443 (99%)	6 (1%)	69	74
All	All	889/940 (95%)	877 (99%)	12 (1%)	69	74

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	65	CYS
1	A	97	PHE
1	A	122	LEU
1	A	125	ASN
1	A	160	MET
1	A	527	GLN
1	B	30	ASP
1	B	65	CYS
1	B	74	LYS
1	B	97	PHE
1	B	125	ASN
1	B	527	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	119	GLN
1	A	282	GLN
1	A	452	GLN
1	A	527	GLN
1	B	58	ASN
1	B	119	GLN
1	B	125	ASN
1	B	527	GLN

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

6 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	HEM	B	601	1	27,50,50	1.77	6 (22%)	17,82,82	1.77	2 (11%)
3	DHE	A	602	1,4	38,56,56	5.68	23 (60%)	37,94,94	4.27	15 (40%)
2	HEM	A	601	1	27,50,50	1.84	6 (22%)	17,82,82	1.44	2 (11%)
3	DHE	B	602	1,4	38,56,56	5.62	23 (60%)	37,94,94	4.22	15 (40%)
4	SO2	A	603	3	2,2,2	0.24	0	1,1,1	0.83	0
4	SO2	B	603	3	2,2,2	0.18	0	1,1,1	0.85	0

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	DHE	B	602	1,4	-	4/12/108/108	-
2	HEM	B	601	1	-	2/6/54/54	-
3	DHE	A	602	1,4	-	5/12/108/108	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	HEM	A	601	1	-	2/6/54/54	-

All (58) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	602	DHE	C4B-NB	-14.42	1.35	1.49
3	B	602	DHE	C4B-NB	-13.84	1.35	1.49
3	B	602	DHE	CHA-C4D	-13.06	1.37	1.51
3	A	602	DHE	C4C-NC	-12.92	1.36	1.49
3	B	602	DHE	C4C-NC	-12.37	1.37	1.49
3	B	602	DHE	CHA-C1A	-11.85	1.39	1.51
3	A	602	DHE	CHA-C4D	-11.73	1.39	1.51
3	A	602	DHE	CHA-C1A	-11.65	1.39	1.51
3	A	602	DHE	C1C-NC	-9.94	1.36	1.49
3	A	602	DHE	CHB-C1B	-9.46	1.37	1.53
3	A	602	DHE	C1B-NB	-9.45	1.36	1.49
3	B	602	DHE	CHB-C1B	-9.22	1.37	1.53
3	B	602	DHE	C1B-NB	-9.17	1.37	1.49
3	B	602	DHE	C1C-NC	-9.02	1.37	1.49
3	A	602	DHE	CHC-C4B	-6.17	1.37	1.52
3	A	602	DHE	CHD-C4C	-6.13	1.37	1.54
3	B	602	DHE	CHC-C4B	-6.00	1.38	1.52
3	B	602	DHE	CHD-C4C	-5.69	1.39	1.54
3	A	602	DHE	CHB-C4A	-4.99	1.39	1.51
3	A	602	DHE	CAB-C3B	-4.97	1.48	1.56
3	B	602	DHE	CHB-C4A	-4.81	1.39	1.51
3	A	602	DHE	CHD-C1D	-4.68	1.39	1.51
3	B	602	DHE	CAC-C3C	-4.57	1.48	1.56
3	B	602	DHE	CHD-C1D	-4.46	1.40	1.51
3	B	602	DHE	CAB-C3B	-4.43	1.49	1.56
3	B	602	DHE	C1D-C2D	4.41	1.44	1.38
3	B	602	DHE	CGC-C3C	4.23	1.61	1.54
3	B	602	DHE	CHC-C1C	-4.18	1.37	1.53
3	A	602	DHE	CAC-C3C	-4.14	1.49	1.56
3	A	602	DHE	CHC-C1C	-4.11	1.37	1.53
3	A	602	DHE	CGB-C3B	4.10	1.61	1.54
2	A	601	HEM	C3C-C2C	-4.04	1.34	1.40
3	B	602	DHE	CGB-C3B	3.82	1.61	1.54
2	B	601	HEM	C3C-C2C	-3.81	1.35	1.40
3	A	602	DHE	C1D-C2D	3.71	1.43	1.38
3	A	602	DHE	C4A-C3A	3.65	1.43	1.38
3	B	602	DHE	C4A-C3A	3.63	1.43	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	601	HEM	CBB-CAB	3.62	1.53	1.29
3	B	602	DHE	CBD-CAD	-3.62	1.28	1.53
2	B	601	HEM	CBC-CAC	3.59	1.53	1.29
3	A	602	DHE	CGC-C3C	3.56	1.60	1.54
2	A	601	HEM	CBC-CAC	3.53	1.52	1.29
2	A	601	HEM	CBB-CAB	3.52	1.52	1.29
3	A	602	DHE	CBD-CAD	-3.41	1.29	1.53
3	B	602	DHE	CAD-C3D	-3.33	1.47	1.52
3	A	602	DHE	C1A-C2A	3.27	1.42	1.38
3	A	602	DHE	CAD-C3D	-3.25	1.47	1.52
2	A	601	HEM	C3B-CAB	3.12	1.54	1.47
2	B	601	HEM	C3C-CAC	3.10	1.54	1.47
3	A	602	DHE	C4D-C3D	3.09	1.42	1.38
2	A	601	HEM	C3B-C2B	-3.01	1.36	1.40
3	B	602	DHE	C4D-C3D	3.00	1.42	1.38
3	B	602	DHE	C1A-C2A	2.97	1.42	1.38
2	B	601	HEM	C3B-CAB	2.68	1.53	1.47
2	A	601	HEM	C3C-CAC	2.65	1.53	1.47
2	B	601	HEM	C3B-C2B	-2.34	1.37	1.40
3	B	602	DHE	C3B-C4B	-2.09	1.52	1.55
3	A	602	DHE	C3C-C4C	-2.01	1.52	1.55

All (34) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	602	DHE	CHC-C1C-NC	12.75	126.79	110.94
3	B	602	DHE	CHC-C1C-NC	12.68	126.71	110.94
3	A	602	DHE	CHB-C1B-NB	8.86	127.46	110.75
3	B	602	DHE	CHB-C1B-NB	8.71	127.18	110.75
3	B	602	DHE	C3B-C4B-NB	8.30	111.97	104.67
3	A	602	DHE	C3C-C4C-NC	8.06	111.76	104.67
3	A	602	DHE	CAD-CBD-CGD	8.00	126.09	112.67
3	B	602	DHE	C3C-C4C-NC	7.74	111.48	104.67
3	A	602	DHE	CBD-CAD-C3D	7.66	126.61	112.49
3	B	602	DHE	CBD-CAD-C3D	7.56	126.43	112.49
3	B	602	DHE	CAD-CBD-CGD	7.48	125.22	112.67
3	A	602	DHE	C3B-C4B-NB	7.45	111.23	104.67
3	A	602	DHE	C4D-CHA-C1A	6.70	129.22	112.87
3	B	602	DHE	C4D-CHA-C1A	6.54	128.82	112.87
2	B	601	HEM	CBD-CAD-C3D	5.87	123.29	112.48
3	A	602	DHE	CAA-C2A-C1A	-5.67	123.31	127.30
3	B	602	DHE	CAD-C3D-C4D	-5.17	123.67	127.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	601	HEM	CBD-CAD-C3D	4.91	121.53	112.48
3	A	602	DHE	CAD-C3D-C4D	-4.84	123.89	127.30
3	B	602	DHE	C1C-CHC-C4B	4.53	129.47	116.16
3	A	602	DHE	C1C-CHC-C4B	4.51	129.40	116.16
3	B	602	DHE	CAA-C2A-C1A	-4.19	124.35	127.30
3	A	602	DHE	CHC-C4B-C3B	3.72	125.64	118.34
3	A	602	DHE	CAA-CBA-CGA	3.65	118.80	112.67
3	B	602	DHE	CAA-CBA-CGA	3.62	118.74	112.67
3	B	602	DHE	CHC-C4B-C3B	3.44	125.08	118.34
3	B	602	DHE	CHB-C1B-C2B	3.08	123.55	114.47
3	A	602	DHE	CHB-C1B-C2B	2.95	123.16	114.47
3	A	602	DHE	CHC-C1C-C2C	2.93	123.84	114.70
3	B	602	DHE	CHC-C1C-C2C	2.84	123.58	114.70
2	A	601	HEM	C3B-C4B-NB	2.38	112.29	109.21
3	A	602	DHE	CMA-C3A-C2A	2.33	129.34	124.94
3	B	602	DHE	CMD-C2D-C3D	2.16	129.02	124.94
2	B	601	HEM	CAD-C3D-C2D	-2.15	121.06	127.25

There are no chirality outliers.

All (13) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	601	HEM	C2D-C3D-CAD-CBD
2	B	601	HEM	C4D-C3D-CAD-CBD
3	A	602	DHE	C2D-C3D-CAD-CBD
3	A	602	DHE	C4D-C3D-CAD-CBD
2	A	601	HEM	C2D-C3D-CAD-CBD
3	B	602	DHE	C2D-C3D-CAD-CBD
3	B	602	DHE	C4D-C3D-CAD-CBD
3	A	602	DHE	C4C-C3C-CAC-CBC
2	A	601	HEM	C4D-C3D-CAD-CBD
3	A	602	DHE	C4B-C3B-CAB-CBB
3	A	602	DHE	CGB-C3B-CAB-CBB
3	B	602	DHE	CGC-C3C-CAC-CBC
3	B	602	DHE	CGB-C3B-CAB-CBB

There are no ring outliers.

4 monomers are involved in 5 short contacts:

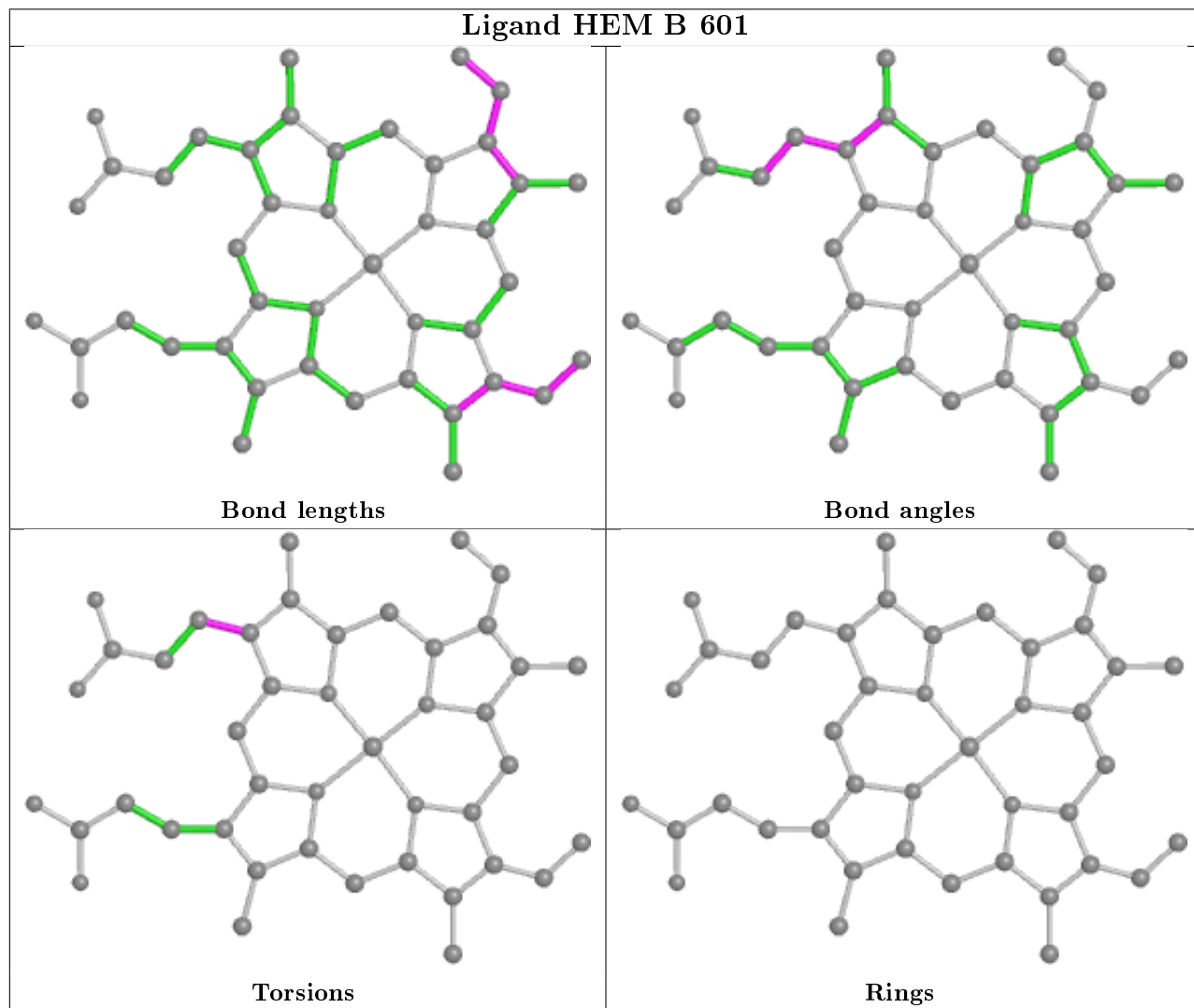
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	601	HEM	1	0

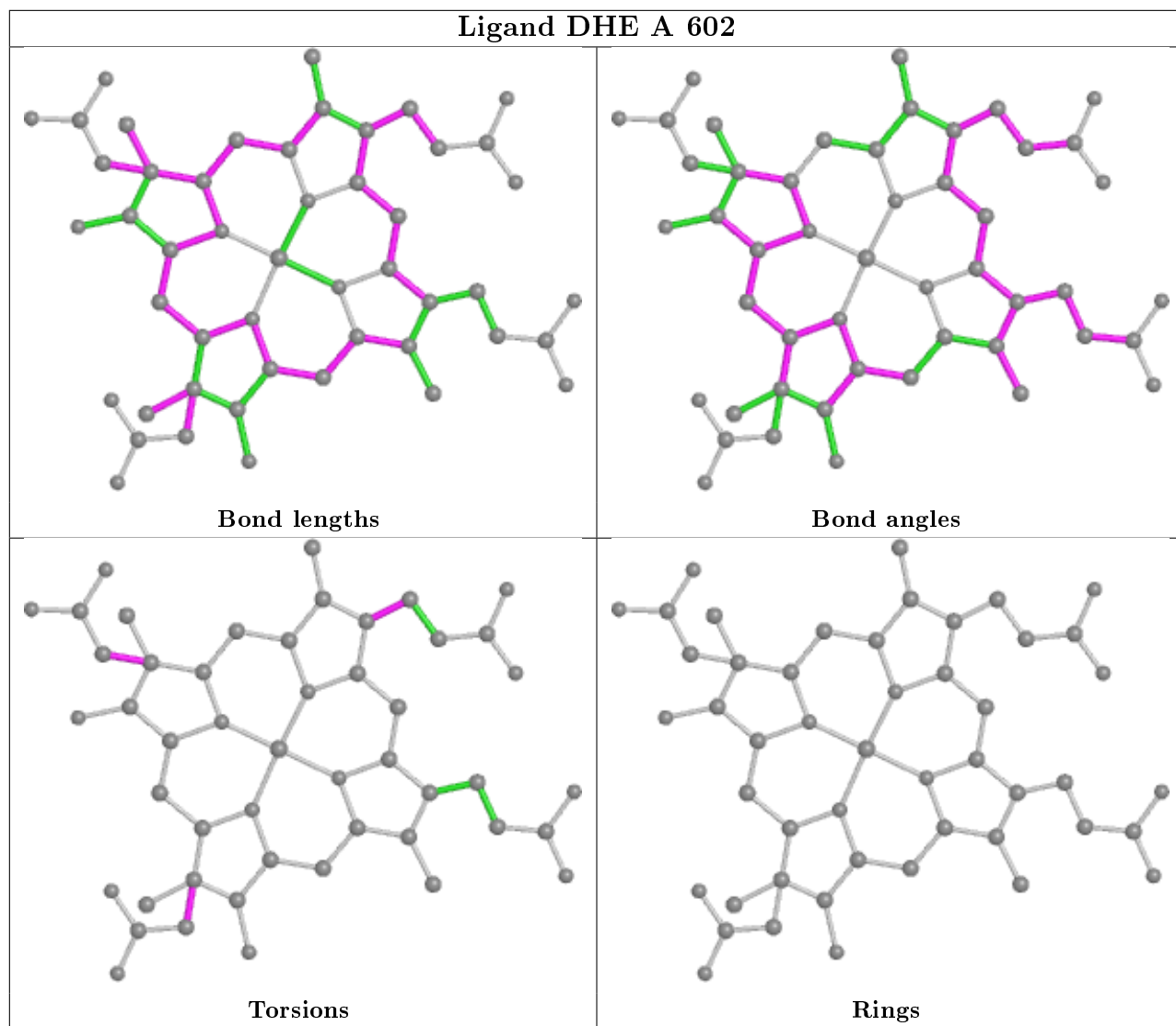
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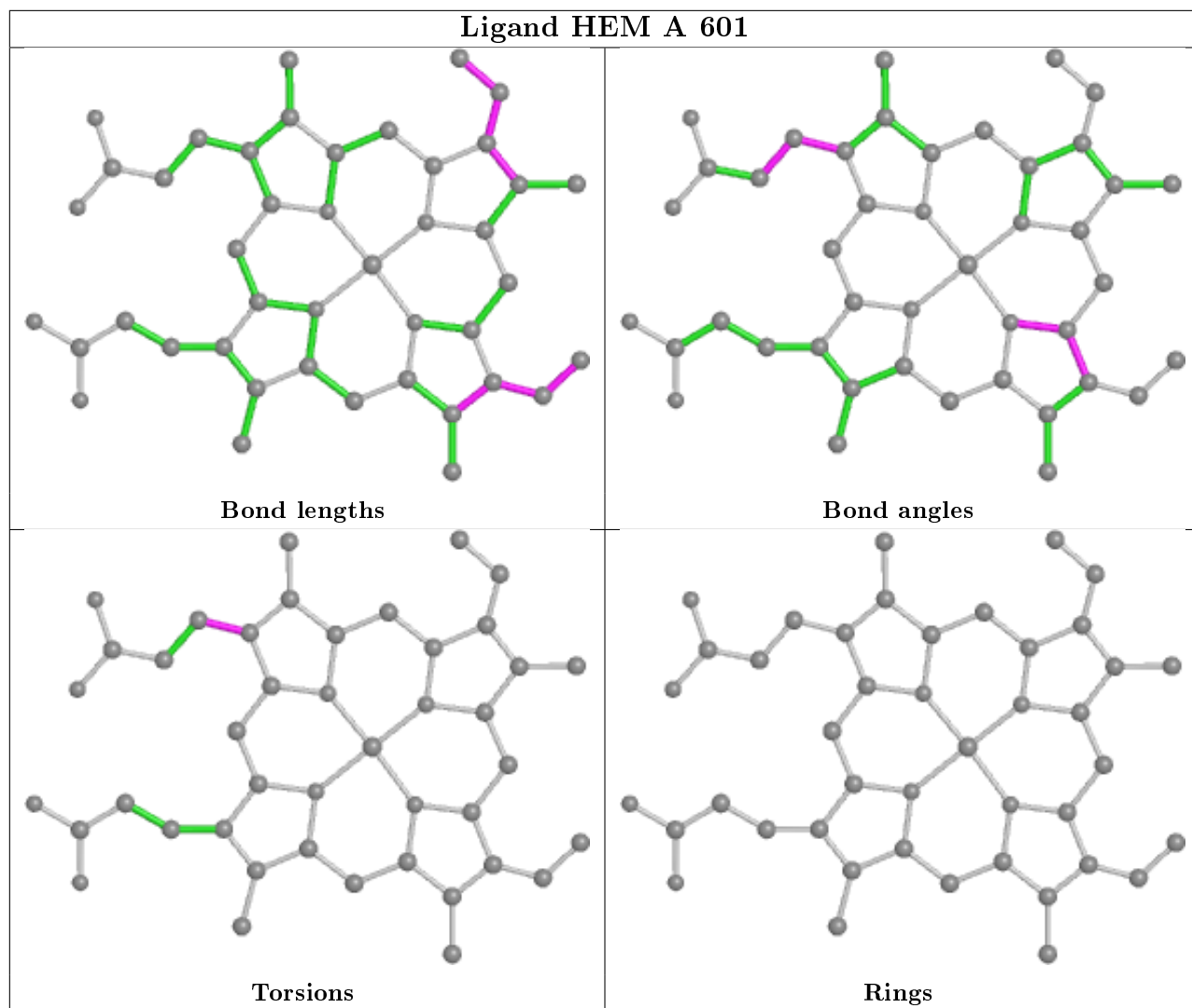
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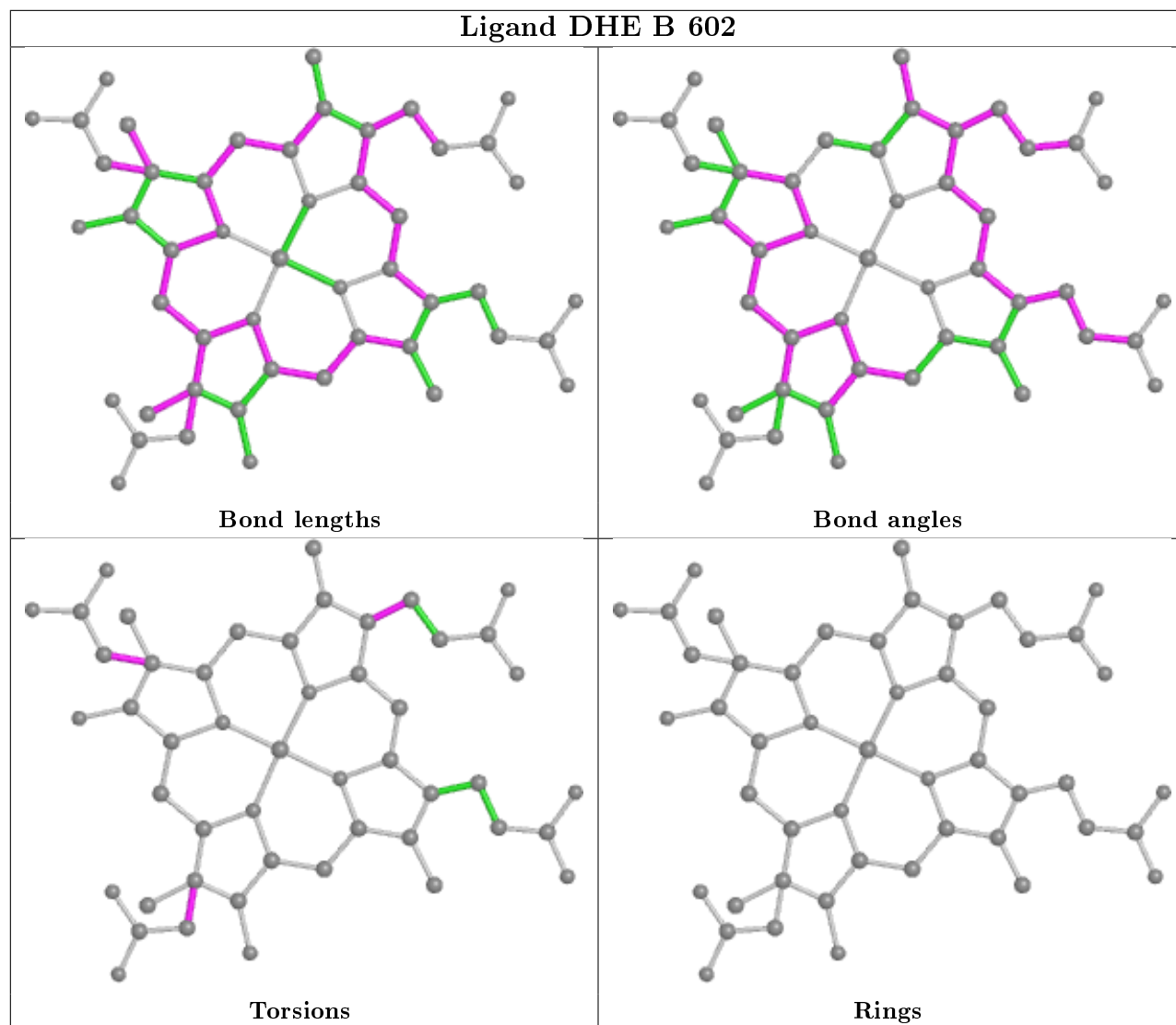
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	602	DHE	1	0
2	A	601	HEM	2	0
3	B	602	DHE	1	0

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 than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.









5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

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	532/567 (93%)	-0.12	45 (8%) 10 10	13, 23, 44, 52	59 (11%)
1	B	542/567 (95%)	-0.39	17 (3%) 49 48	14, 24, 44, 52	0
All	All	1074/1134 (94%)	-0.26	62 (5%) 23 22	13, 23, 44, 52	59 (5%)

All (62) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	114	GLU	9.6
1	A	105	GLY	9.3
1	A	103	PRO	8.2
1	A	98	ILE	7.0
1	A	108	ASN	6.5
1	B	103	PRO	6.4
1	A	100	TYR	6.2
1	A	62	PHE	5.7
1	A	111	THR	5.4
1	A	97	PHE	5.3
1	A	113	GLY	5.0
1	B	499	THR	4.9
1	A	110	GLY	4.8
1	A	93	TYR	4.7
1	A	102	SER	4.5
1	A	94	LEU	4.4
1	A	81	LEU	4.4
1	B	104	ALA	4.4
1	A	71	VAL	4.3
1	A	109	TRP	4.2
1	A	96	SER	3.9
1	A	104	ALA	3.9
1	A	106	MET	3.9
1	A	112	SER	3.8

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Mol	Chain	Res	Type	RSRZ
1	A	120	VAL	3.7
1	A	115	LEU	3.7
1	A	99	THR	3.7
1	B	502	GLN	3.4
1	B	34	GLN	3.4
1	A	64	ARG	3.3
1	B	105	GLY	3.2
1	A	107	PRO	3.1
1	A	72	LEU	3.0
1	A	501	GLY	3.0
1	A	101	ALA	2.9
1	A	526	ASP	2.8
1	A	82	THR	2.8
1	A	197	TYR	2.8
1	A	36	ASP	2.8
1	A	133	ALA	2.8
1	A	65	CYS	2.8
1	A	92	ASP	2.7
1	B	36	ASP	2.7
1	A	134	PRO	2.7
1	B	35	GLN	2.7
1	A	121	ASP	2.6
1	A	44	GLU	2.6
1	A	160	MET	2.5
1	A	95	GLN	2.4
1	A	66	ALA	2.4
1	B	526	ASP	2.4
1	A	73	ARG	2.3
1	B	160	MET	2.3
1	B	482	GLY	2.3
1	B	101	ALA	2.2
1	A	69	HIS	2.2
1	B	500	GLU	2.2
1	B	483	SER	2.2
1	A	68	CYS	2.1
1	B	462	PRO	2.1
1	B	196	GLY	2.1
1	B	52	ALA	2.0

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 [i](#)

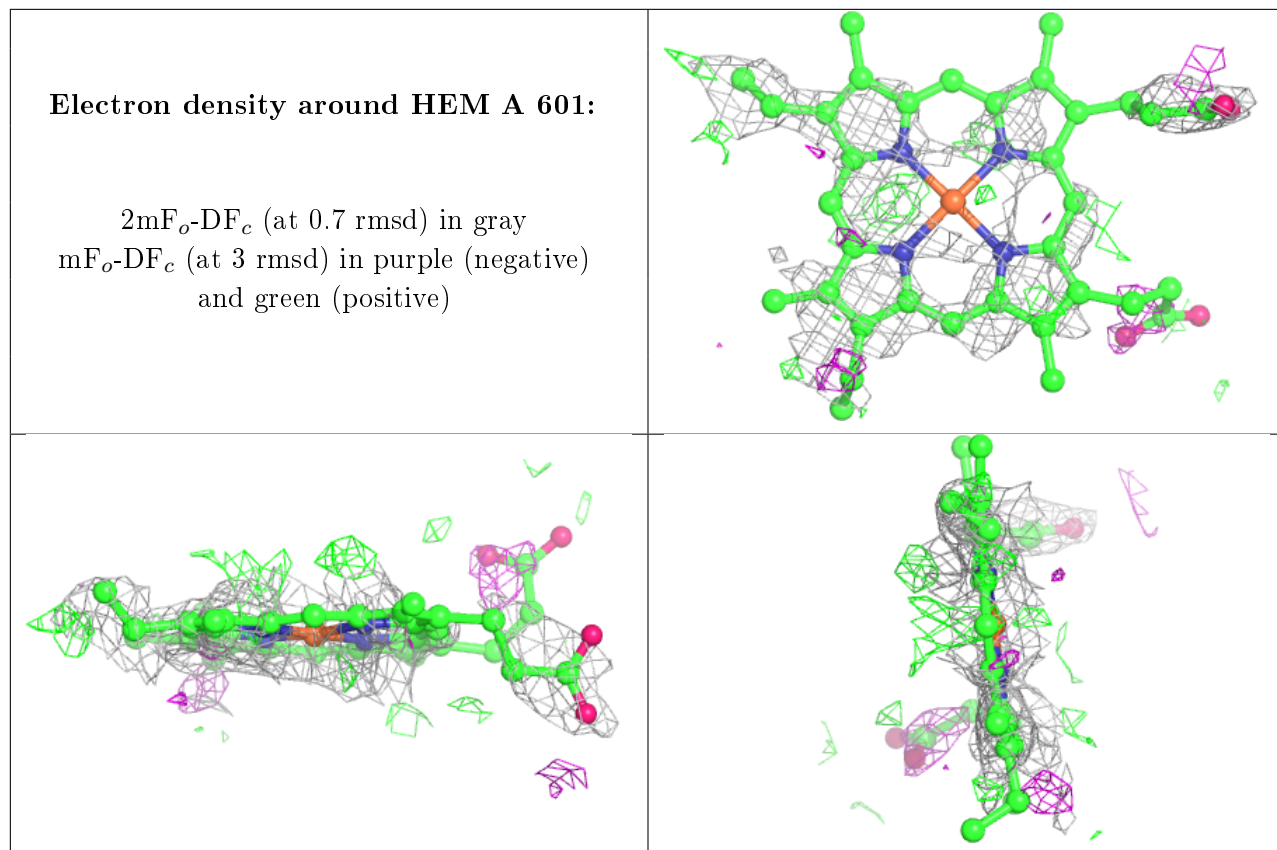
There are no carbohydrates in this entry.

6.4 Ligands [i](#)

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

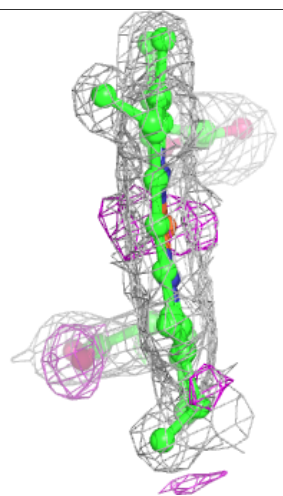
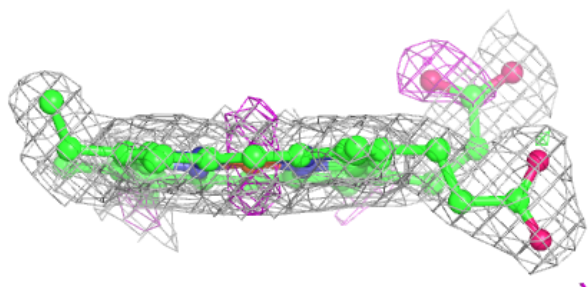
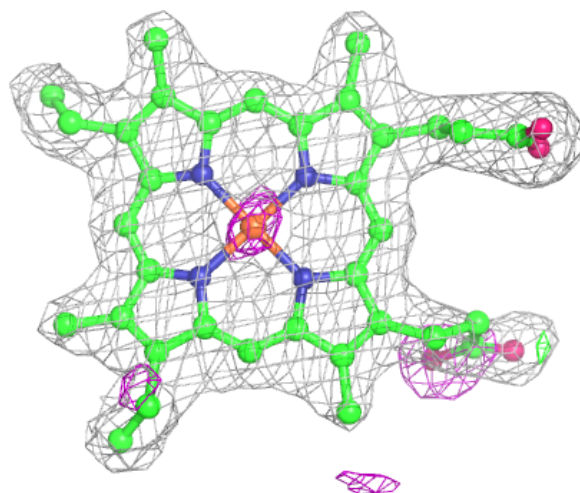
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	HEM	A	601	43/43	0.71	0.37	30,37,45,48	43
2	HEM	B	601	43/43	0.92	0.14	26,31,45,52	0
3	DHE	A	602	49/49	0.98	0.08	12,17,26,33	0
4	SO2	B	603	3/3	0.98	0.08	23,23,30,31	0
4	SO2	A	603	3/3	0.99	0.11	24,24,26,30	0
3	DHE	B	602	49/49	0.99	0.06	14,18,23,28	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.



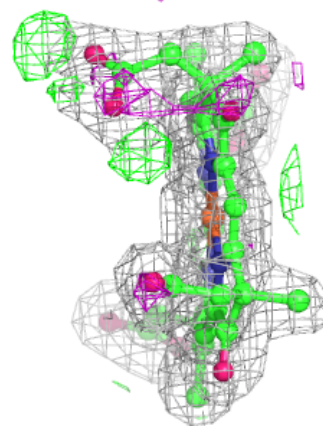
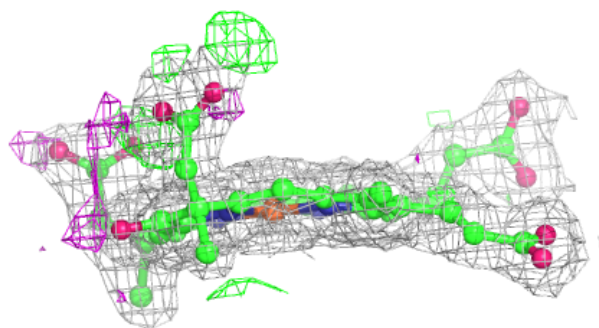
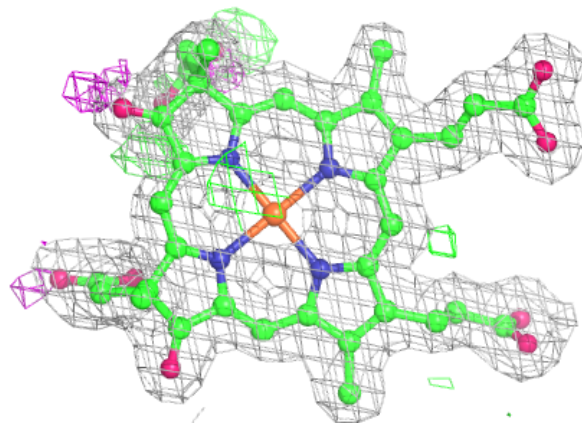
Electron density around HEM B 601:

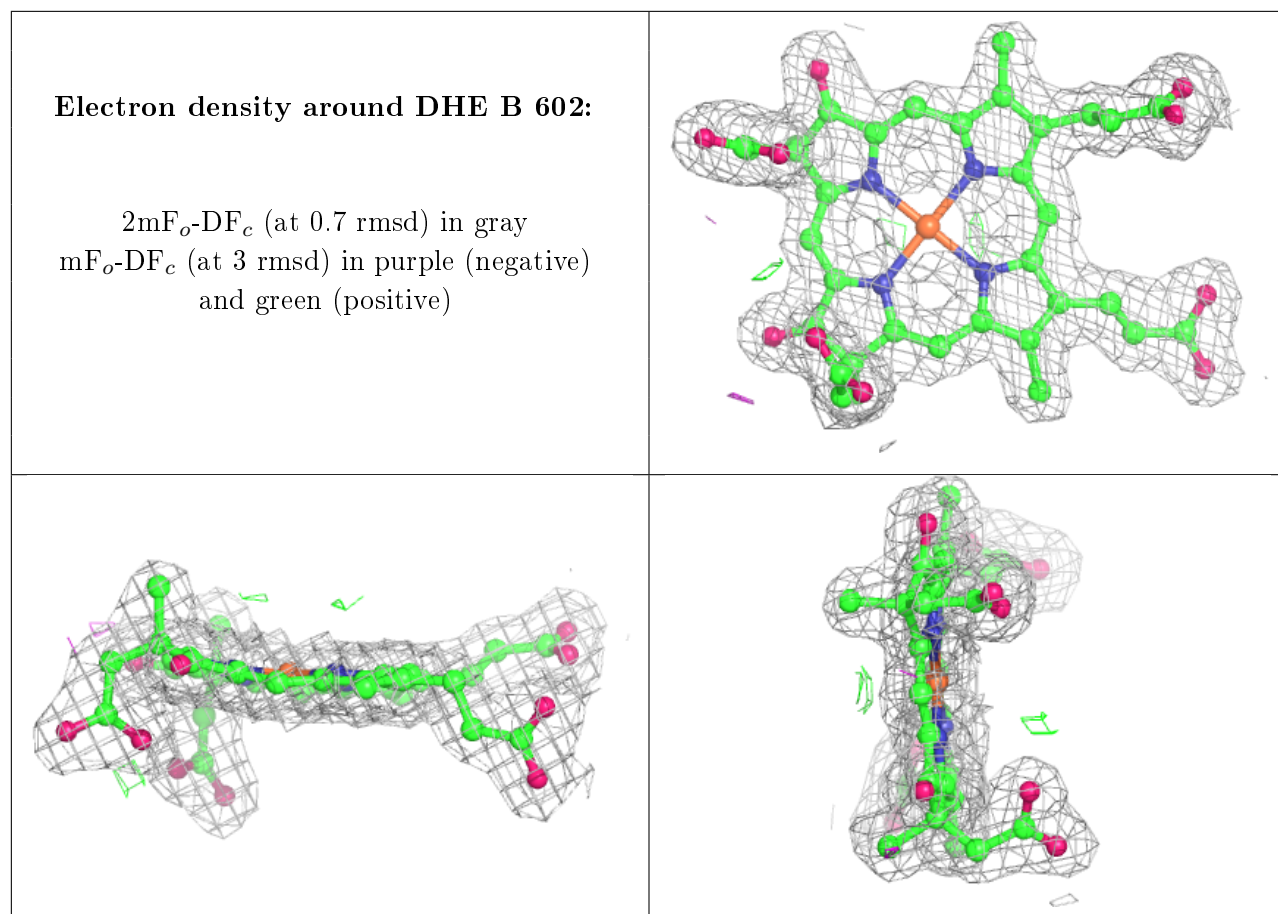
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around DHE A 602:

$2mF_o-DF_c$ (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.