



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 26, 2026 – 01:36 PM EDT

PDB ID : 8FDK / pdb_00008fdk
Title : Phenylhydroxylamine in Reaction with Human Hemoglobin
Authors : Powell, S.M.; Thomas, L.M.; Richter-Addo, G.B.
Deposited on : 2022-12-03
Resolution : 1.89 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

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	GOL	C	203	-	-	X	-

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 4779 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.

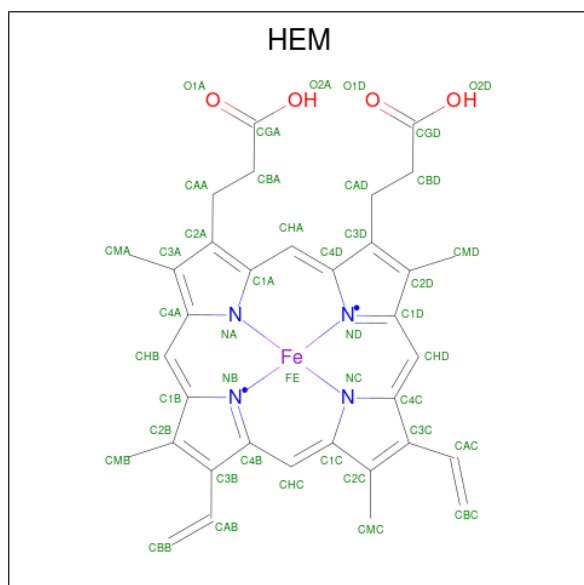
- Molecule 1 is a protein called Hemoglobin subunit alpha.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	141	Total	C	N	O	S	0	0	0
			1068	685	187	193	3			
1	C	141	Total	C	N	O	S	0	0	0
			1068	685	187	193	3			

- Molecule 2 is a protein called Hemoglobin subunit beta.

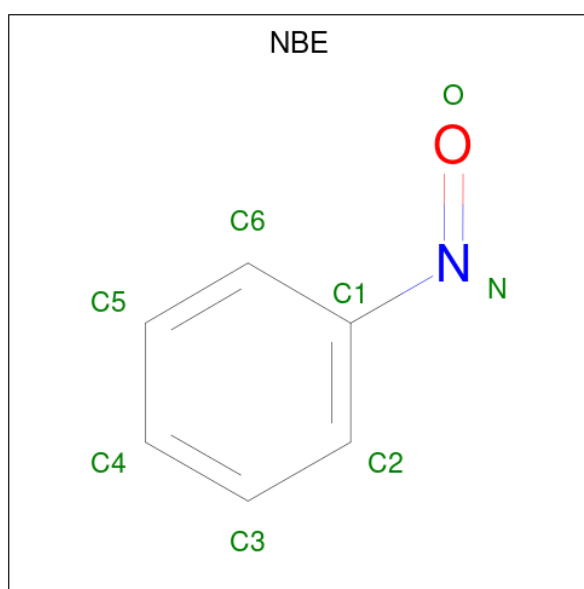
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	143	Total	C	N	O	S	0	3	0
			1120	718	194	205	3			
2	D	143	Total	C	N	O	S	0	2	0
			1112	714	193	202	3			

- Molecule 3 is PROTOPORPHYRIN IX CONTAINING FE (CCD ID: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).



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

- Molecule 4 is NITROSOBENZENE (CCD ID: NBE) (formula: C_6H_5NO) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	B	1	Total 8	C 6	N 1	O 1	0	0
4	B	1	Total 8	C 6	N 1	O 1	0	0
4	C	1	Total 8	C 6	N 1	O 1	0	0
4	D	1	Total 8	C 6	N 1	O 1	0	0
4	D	1	Total 8	C 6	N 1	O 1	0	0

- Molecule 5 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



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

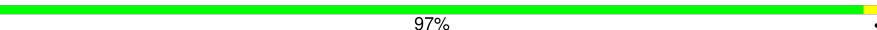
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	62	Total	O	0	0
			62	62		
6	B	34	Total	O	0	0
			34	34		
6	C	64	Total	O	0	0
			64	64		
6	D	33	Total	O	0	0
			33	33		

3 Residue-property plots [i](#)

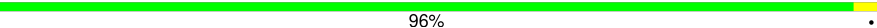
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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.

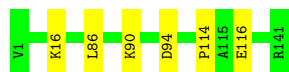
- Molecule 1: Hemoglobin subunit alpha

Chain A:  97% .



- Molecule 1: Hemoglobin subunit alpha

Chain C:  96% .



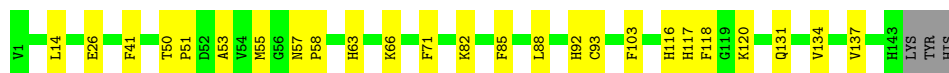
- Molecule 2: Hemoglobin subunit beta

Chain B:  84% 14% .



- Molecule 2: Hemoglobin subunit beta

Chain D:  81% 17% .



4 Data and refinement statistics

Property	Value	Source
Space group	P 41	Depositor
Cell constants a, b, c, α , β , γ	53.26Å 53.26Å 194.71Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	32.58 – 1.89 32.58 – 1.89	Depositor EDS
% Data completeness (in resolution range)	99.5 (32.58-1.89) 99.7 (32.58-1.89)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.80 (at 1.88Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.208 , 0.246 0.208 , 0.248	Depositor DCC
R_{free} test set	2135 reflections (4.91%)	wwPDB-VP
Wilson B-factor (Å ²)	31.3	Xtriage
Anisotropy	0.481	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 36.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.450 for h,-k,-l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	4779	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.24% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: NBE, HEM, GOL, CSD

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	0.32	0/1096	0.50	0/1491
1	C	0.33	0/1096	0.48	0/1491
2	B	0.27	0/1141	0.49	0/1552
2	D	0.28	0/1133	0.46	0/1541
All	All	0.30	0/4466	0.48	0/6075

There are no bond length outliers.

There are no bond angle outliers.

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	1068	0	1073	3	0
1	C	1068	0	1073	4	0
2	B	1120	0	1102	20	0
2	D	1112	0	1099	16	0
3	A	43	0	30	1	0
3	B	43	0	30	3	0
3	C	43	0	30	0	0
3	D	43	0	30	3	0
4	B	16	0	10	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	C	8	0	5	0	0
4	D	16	0	10	2	0
5	C	6	0	8	4	0
6	A	62	0	0	1	0
6	B	34	0	0	0	0
6	C	64	0	0	1	0
6	D	33	0	0	2	0
All	All	4779	0	4500	44	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 (44) 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)
3:B:201:HEM:HBB2	3:B:201:HEM:HHC	1.71	0.72
2:B:98:VAL:HG12	5:C:203:GOL:H32	1.76	0.68
3:D:201:HEM:HBB2	3:D:201:HEM:HHC	1.75	0.67
2:B:88:LEU:HD12	2:B:92:HIS:HE1	1.66	0.60
6:A:308:HOH:O	2:B:120:LYS:HD2	2.01	0.60
2:B:99:ASP:N	5:C:203:GOL:O1	2.37	0.58
6:C:303:HOH:O	2:D:120:LYS:HD2	2.04	0.58
2:B:92:HIS:HB3	2:B:95:LYS:HB2	1.86	0.57
2:B:66:LYS:NZ	3:B:201:HEM:O2A	2.25	0.54
2:D:26[B]:GLU:OE1	2:D:117[B]:HIS:NE2	2.32	0.54
3:D:201:HEM:HMC2	3:D:201:HEM:HBC2	1.90	0.53
2:B:100:PRO:HB2	2:B:104:ARG:NH2	2.23	0.53
2:D:51:PRO:O	2:D:55:MET:HG2	2.08	0.53
2:B:88:LEU:HD12	2:B:92:HIS:CE1	2.45	0.52
1:C:16:LYS:HG3	1:C:116:GLU:HG2	1.93	0.50
2:D:63:HIS:HA	2:D:66:LYS:HG3	1.93	0.49
2:D:88:LEU:O	2:D:92:HIS:ND1	2.38	0.48
2:B:98:VAL:HB	5:C:203:GOL:H12	1.95	0.48
2:B:50:THR:H	2:B:53:ALA:HB3	1.79	0.48
1:C:114:PRO:HA	2:D:116:HIS:CD2	2.50	0.47
2:D:63:HIS:HE1	6:D:303:HOH:O	1.97	0.47
2:B:71:PHE:HD2	4:B:202:NBE:O	1.99	0.46
2:D:103:PHE:CZ	4:D:203:NBE:H6	2.51	0.46
1:C:94:ASP:OD2	5:C:203:GOL:H11	2.16	0.46
1:C:86:LEU:HD12	1:C:90:LYS:HD3	1.99	0.45
2:B:95:LYS:HE3	3:B:201:HEM:CBC	2.46	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:82:LYS:NZ	6:D:302:HOH:O	2.47	0.45
1:A:20:HIS:HB3	1:A:24:TYR:CE1	2.52	0.44
2:B:103:PHE:CZ	4:B:203:NBE:H2	2.52	0.44
2:B:57:ASN:O	2:B:61:LYS:HG3	2.17	0.44
2:B:88:LEU:O	2:B:92:HIS:ND1	2.45	0.44
2:B:88:LEU:CD1	2:B:92:HIS:HE1	2.30	0.44
2:B:61:LYS:HB3	2:B:61:LYS:HE2	1.46	0.44
1:A:61:LYS:HD2	3:A:201:HEM:O1A	2.18	0.43
2:D:85:PHE:CE2	2:D:137:VAL:HG13	2.55	0.42
2:D:41:PHE:O	3:D:201:HEM:HBC2	2.20	0.42
2:D:71:PHE:HD2	4:D:202:NBE:O	2.03	0.41
1:A:114:PRO:HA	2:B:116:HIS:CD2	2.55	0.41
2:B:82:LYS:HA	2:B:82:LYS:HD2	1.76	0.41
2:D:50:THR:CG2	2:D:53:ALA:H	2.33	0.41
2:D:14:LEU:HD21	2:D:118:PHE:CD2	2.56	0.41
2:B:24:GLY:CA	2:B:68:LEU:HD13	2.51	0.41
2:D:131:GLN:HA	2:D:134:VAL:HG22	2.02	0.40
2:D:57:ASN:HA	2:D:58:PRO:HD3	1.95	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	139/141 (99%)	138 (99%)	1 (1%)	0	100	100
1	C	139/141 (99%)	138 (99%)	1 (1%)	0	100	100
2	B	143/146 (98%)	139 (97%)	4 (3%)	0	100	100
2	D	142/146 (97%)	139 (98%)	3 (2%)	0	100	100
All	All	563/574 (98%)	554 (98%)	9 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

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	113/113 (100%)	113 (100%)	0	100	100
1	C	113/113 (100%)	113 (100%)	0	100	100
2	B	117/117 (100%)	117 (100%)	0	100	100
2	D	116/117 (99%)	116 (100%)	0	100	100
All	All	459/460 (100%)	459 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	97	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

2 non-standard protein/DNA/RNA residues 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	CSD	B	93	2	4,7,8	0.94	0	1,8,10	2.26	1 (100%)
2	CSD	D	93	2	4,7,8	1.39	1 (25%)	1,8,10	0.95	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
2	CSD	B	93	2	-	0/2/6/8	-
2	CSD	D	93	2	-	1/2/6/8	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	93	CSD	OD1-SG	-2.41	1.45	1.47

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	93	CSD	OD1-SG-CB	-2.26	101.44	105.60

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	D	93	CSD	CA-CB-SG-OD1

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

10 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
3	HEM	B	201	2	50,50,50	1.55	9 (18%)	67,82,82	1.36	9 (13%)
5	GOL	C	203	-	5,5,5	0.64	0	5,5,5	1.02	0
4	NBE	B	202	-	8,8,8	0.74	0	9,9,9	0.47	0
4	NBE	B	203	-	8,8,8	0.73	0	9,9,9	0.43	0
4	NBE	D	203	-	8,8,8	0.74	0	9,9,9	0.51	0
4	NBE	C	202	-	8,8,8	0.73	0	9,9,9	0.34	0
3	HEM	D	201	2	50,50,50	1.55	10 (20%)	67,82,82	1.50	9 (13%)
3	HEM	A	201	6,1	50,50,50	1.35	8 (16%)	67,82,82	1.25	7 (10%)
4	NBE	D	202	-	8,8,8	0.75	0	9,9,9	0.52	0
3	HEM	C	201	6,1	50,50,50	1.45	11 (22%)	67,82,82	1.31	7 (10%)

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	HEM	B	201	2	-	4/14/54/54	-
5	GOL	C	203	-	-	4/4/4/4	-
4	NBE	B	202	-	-	2/2/2/2	0/1/1/1
4	NBE	B	203	-	-	2/2/2/2	0/1/1/1
4	NBE	D	203	-	-	0/2/2/2	0/1/1/1
4	NBE	C	202	-	-	2/2/2/2	0/1/1/1
3	HEM	D	201	2	-	7/14/54/54	-
3	HEM	A	201	6,1	-	2/14/54/54	-
4	NBE	D	202	-	-	2/2/2/2	0/1/1/1
3	HEM	C	201	6,1	-	3/14/54/54	-

All (38) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	201	HEM	FE-ND	4.94	2.10	1.94
3	B	201	HEM	FE-NB	4.48	2.08	1.94
3	C	201	HEM	FE-NB	4.12	2.07	1.94
3	B	201	HEM	FE-ND	3.99	2.07	1.94
3	D	201	HEM	FE-NB	3.81	2.06	1.94
3	A	201	HEM	FE-NB	3.61	2.06	1.94
3	D	201	HEM	FE-NA	3.61	2.07	1.95
3	B	201	HEM	FE-NA	3.24	2.05	1.95

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	201	HEM	FE-ND	3.19	2.04	1.94
3	A	201	HEM	CAB-C3B	3.11	1.55	1.47
3	D	201	HEM	CAC-C3C	3.04	1.55	1.47
3	B	201	HEM	CAB-C3B	3.03	1.55	1.47
3	D	201	HEM	CAB-C3B	3.00	1.55	1.47
3	C	201	HEM	CAB-C3B	2.98	1.55	1.47
3	C	201	HEM	CAC-C3C	2.92	1.55	1.47
3	A	201	HEM	CAC-C3C	2.81	1.54	1.47
3	B	201	HEM	CAC-C3C	2.78	1.54	1.47
3	C	201	HEM	FE-NA	2.72	2.04	1.95
3	B	201	HEM	FE-NC	2.61	2.03	1.95
3	D	201	HEM	CMA-C3A	2.57	1.56	1.50
3	A	201	HEM	FE-ND	2.49	2.02	1.94
3	A	201	HEM	FE-NC	2.48	2.03	1.95
3	C	201	HEM	FE-NC	2.47	2.03	1.95
3	B	201	HEM	CMA-C3A	2.45	1.55	1.50
3	A	201	HEM	CMD-C2D	2.42	1.55	1.50
3	A	201	HEM	CMA-C3A	2.37	1.55	1.50
3	C	201	HEM	CMD-C2D	2.32	1.55	1.50
3	A	201	HEM	CMB-C2B	2.22	1.55	1.50
3	D	201	HEM	CMD-C2D	2.17	1.55	1.50
3	C	201	HEM	CMA-C3A	2.17	1.55	1.50
3	D	201	HEM	CMB-C2B	2.15	1.55	1.50
3	C	201	HEM	CMC-C2C	2.13	1.55	1.50
3	C	201	HEM	CMB-C2B	2.13	1.55	1.50
3	B	201	HEM	CMD-C2D	2.13	1.55	1.50
3	C	201	HEM	C2A-C3A	-2.10	1.33	1.38
3	B	201	HEM	CMB-C2B	2.08	1.55	1.50
3	D	201	HEM	CAA-C2A	2.08	1.56	1.51
3	D	201	HEM	CMC-C2C	2.07	1.55	1.50

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	201	HEM	C1B-NB-C4B	3.52	109.37	105.21
3	D	201	HEM	C3B-C4B-NB	-3.50	106.95	109.47
3	B	201	HEM	C4D-ND-C1D	3.38	109.20	105.21
3	C	201	HEM	CHC-C4B-NB	3.26	127.93	124.42
3	A	201	HEM	CHC-C4B-NB	3.10	127.76	124.42
3	C	201	HEM	CBD-CAD-C3D	-2.97	104.33	112.53
3	A	201	HEM	CBD-CAD-C3D	-2.93	104.44	112.53
3	D	201	HEM	CHD-C4C-NC	2.87	127.58	124.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	201	HEM	C2A-C1A-NA	-2.85	106.99	110.15
3	C	201	HEM	C4D-ND-C1D	2.82	108.54	105.21
3	B	201	HEM	C1B-NB-C4B	2.80	108.52	105.21
3	D	201	HEM	CAA-C2A-C3A	2.76	133.27	127.07
3	B	201	HEM	C3B-C2B-C1B	2.76	108.48	106.41
3	D	201	HEM	C3B-C2B-C1B	2.74	108.47	106.41
3	D	201	HEM	CAA-C2A-C1A	-2.74	119.60	124.94
3	D	201	HEM	CBA-CAA-C2A	2.72	120.05	112.53
3	B	201	HEM	C2D-C1D-ND	-2.66	106.83	109.90
3	B	201	HEM	CHC-C1C-NC	2.66	127.35	124.45
3	D	201	HEM	C2B-C1B-NB	-2.55	106.91	109.84
3	C	201	HEM	C3D-C4D-ND	-2.50	107.43	110.17
3	D	201	HEM	C4D-ND-C1D	2.49	108.16	105.21
3	B	201	HEM	C3B-C4B-NB	-2.44	107.72	109.47
3	A	201	HEM	C2A-C1A-NA	-2.42	107.47	110.15
3	B	201	HEM	C3D-C4D-ND	-2.40	107.53	110.17
3	C	201	HEM	CAA-CBA-CGA	-2.35	107.44	113.67
3	A	201	HEM	C4D-ND-C1D	2.32	107.95	105.21
3	A	201	HEM	C3D-C4D-ND	-2.31	107.64	110.17
3	B	201	HEM	C2A-C1A-NA	-2.09	107.83	110.15
3	A	201	HEM	CHD-C1D-ND	2.06	126.64	124.42
3	A	201	HEM	CAA-CBA-CGA	-2.06	108.21	113.67
3	B	201	HEM	C2B-C1B-NB	-2.05	107.48	109.84
3	C	201	HEM	CHB-C4A-NA	2.02	127.52	123.86

There are no chirality outliers.

All (28) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	D	201	HEM	C3A-C2A-CAA-CBA
3	D	201	HEM	C2D-C3D-CAD-CBD
3	D	201	HEM	C4D-C3D-CAD-CBD
4	B	202	NBE	C6-C1-N-O
4	B	203	NBE	C2-C1-N-O
4	B	203	NBE	C6-C1-N-O
4	C	202	NBE	C2-C1-N-O
4	C	202	NBE	C6-C1-N-O
4	D	202	NBE	C6-C1-N-O
5	C	203	GOL	O1-C1-C2-C3
5	C	203	GOL	C1-C2-C3-O3
5	C	203	GOL	O2-C2-C3-O3
3	D	201	HEM	C1A-C2A-CAA-CBA

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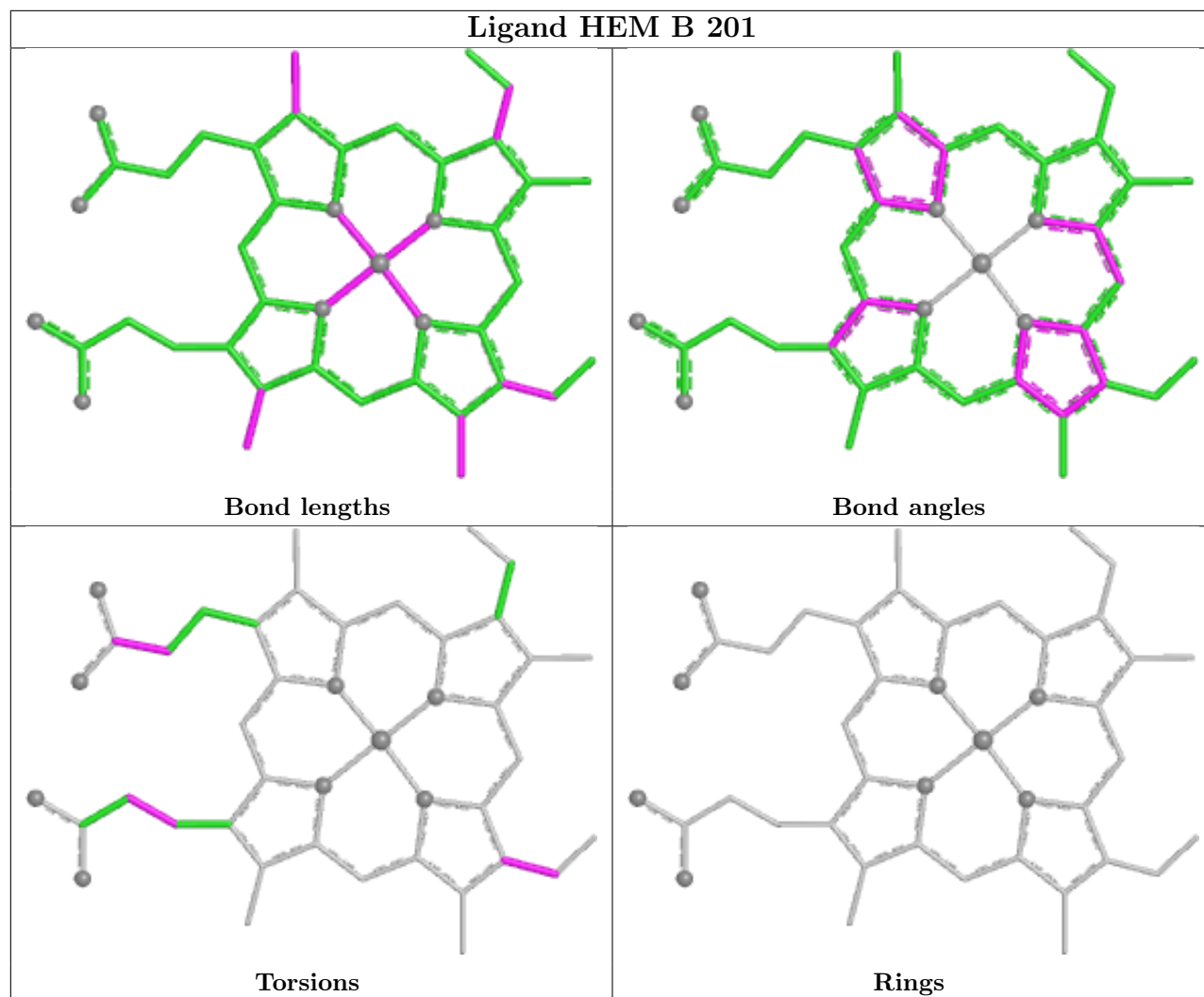
Mol	Chain	Res	Type	Atoms
5	C	203	GOL	O1-C1-C2-O2
4	D	202	NBE	C2-C1-N-O
3	B	201	HEM	C4B-C3B-CAB-CBB
3	D	201	HEM	C4B-C3B-CAB-CBB
3	B	201	HEM	C2A-CAA-CBA-CGA
3	C	201	HEM	CAD-CBD-CGD-O2D
4	B	202	NBE	C2-C1-N-O
3	C	201	HEM	CAD-CBD-CGD-O1D
3	A	201	HEM	CAD-CBD-CGD-O2D
3	A	201	HEM	CAD-CBD-CGD-O1D
3	D	201	HEM	CAD-CBD-CGD-O1D
3	B	201	HEM	CAD-CBD-CGD-O2D
3	B	201	HEM	CAD-CBD-CGD-O1D
3	D	201	HEM	CAD-CBD-CGD-O2D
3	C	201	HEM	CAA-CBA-CGA-O2A

There are no ring outliers.

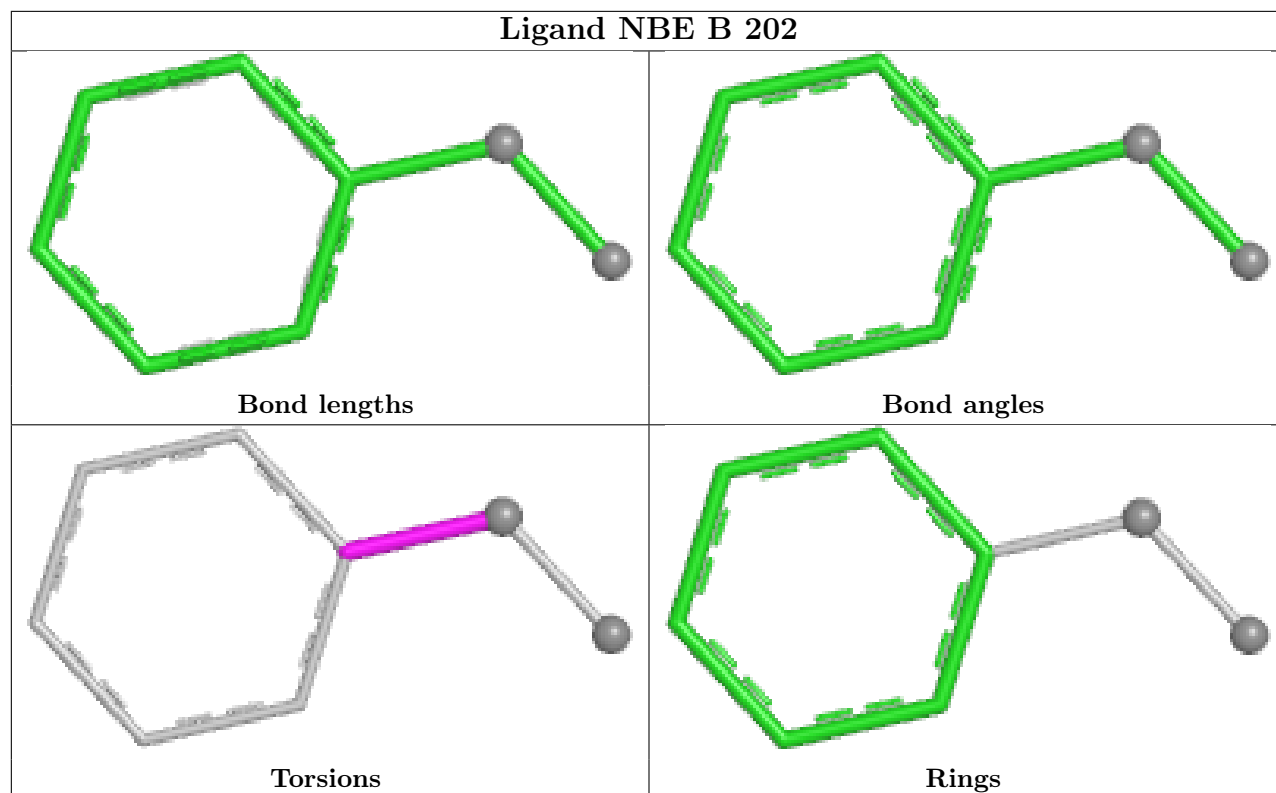
8 monomers are involved in 15 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	201	HEM	3	0
5	C	203	GOL	4	0
4	B	202	NBE	1	0
4	B	203	NBE	1	0
4	D	203	NBE	1	0
3	D	201	HEM	3	0
3	A	201	HEM	1	0
4	D	202	NBE	1	0

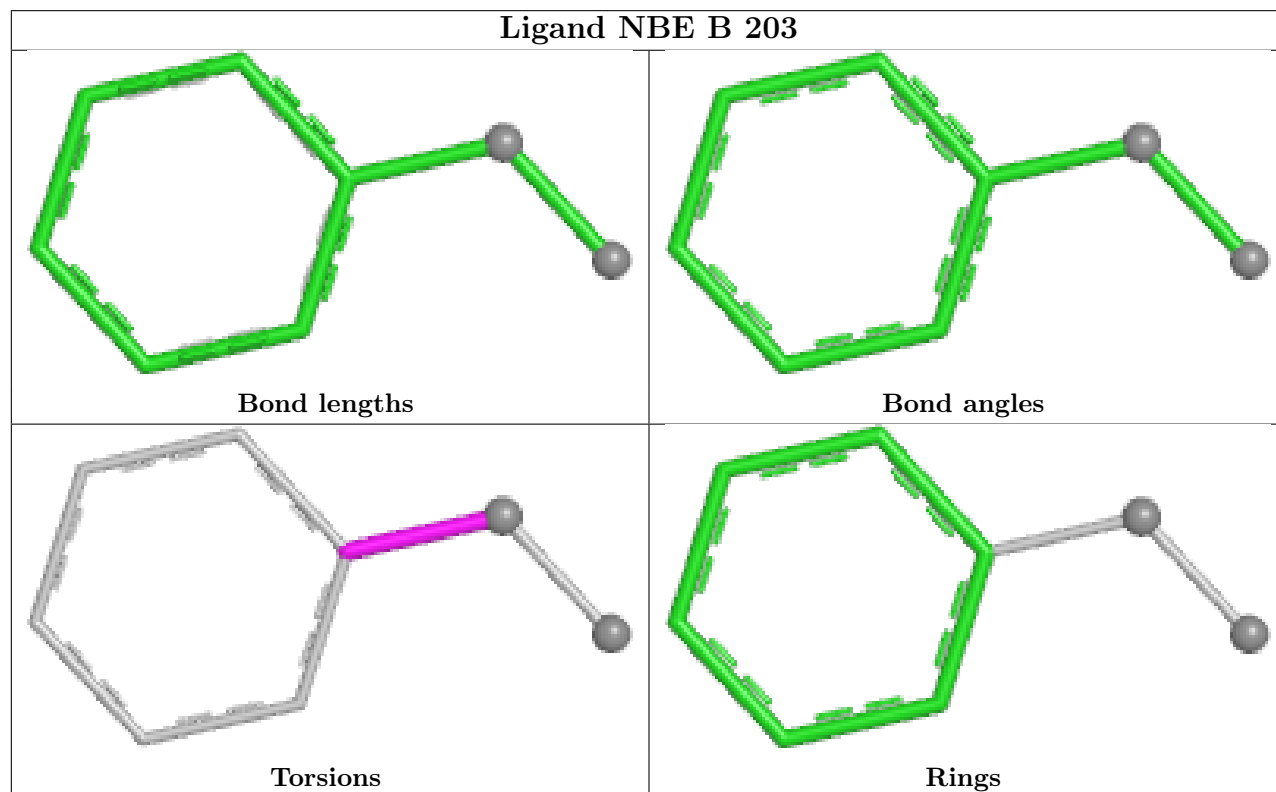
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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.



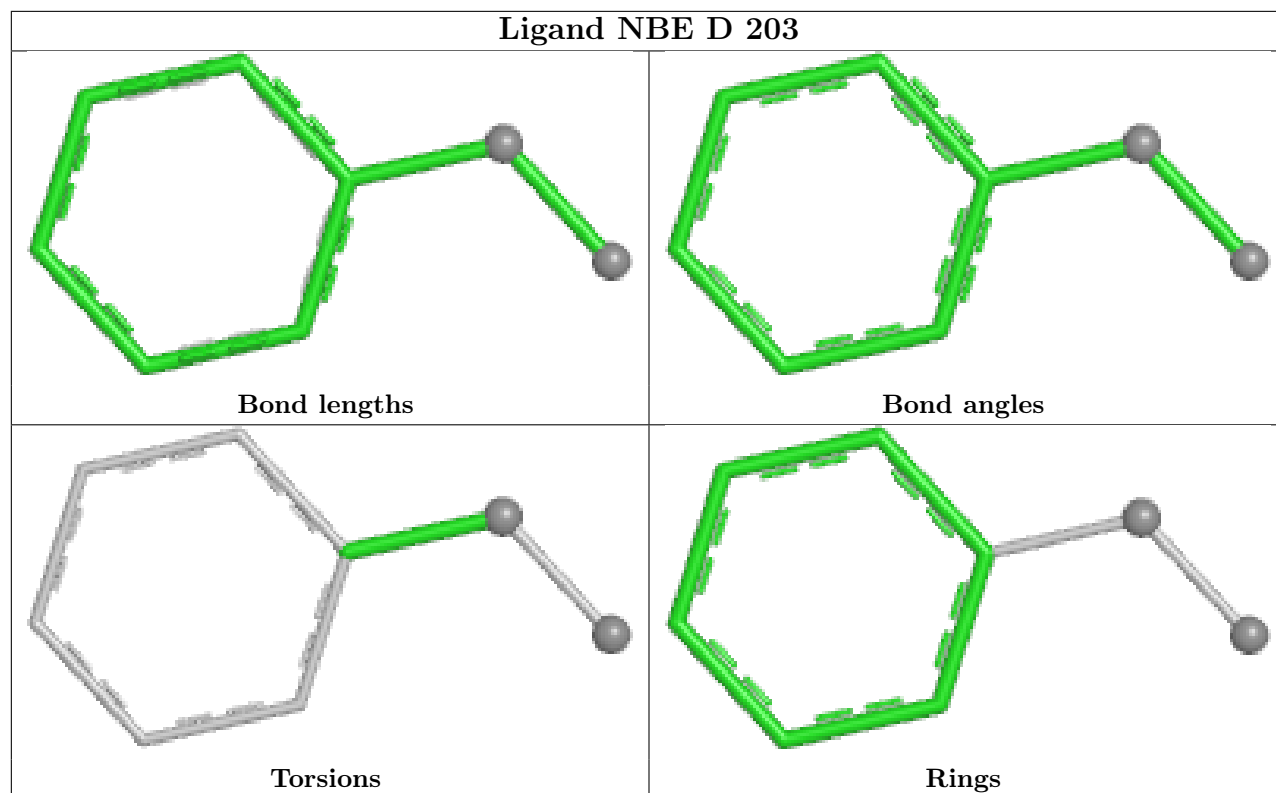
Ligand NBE B 202



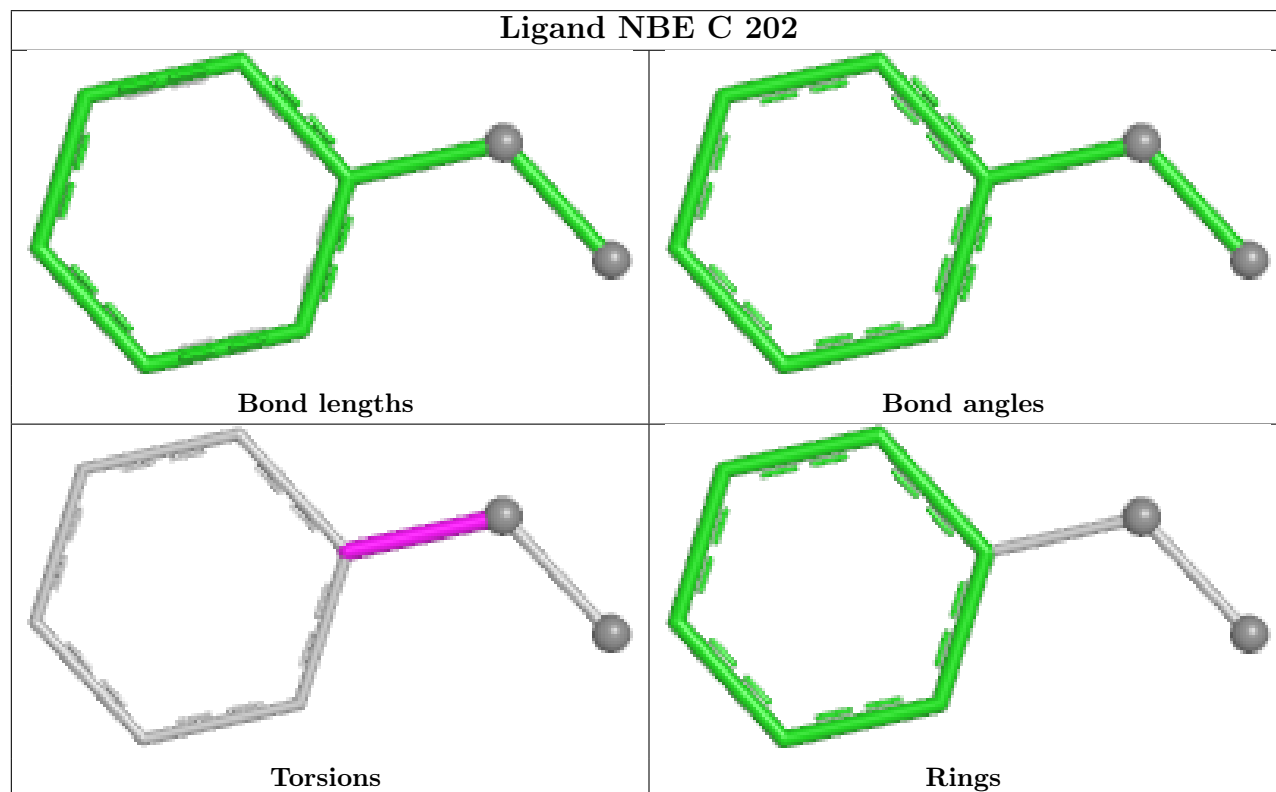
Ligand NBE B 203

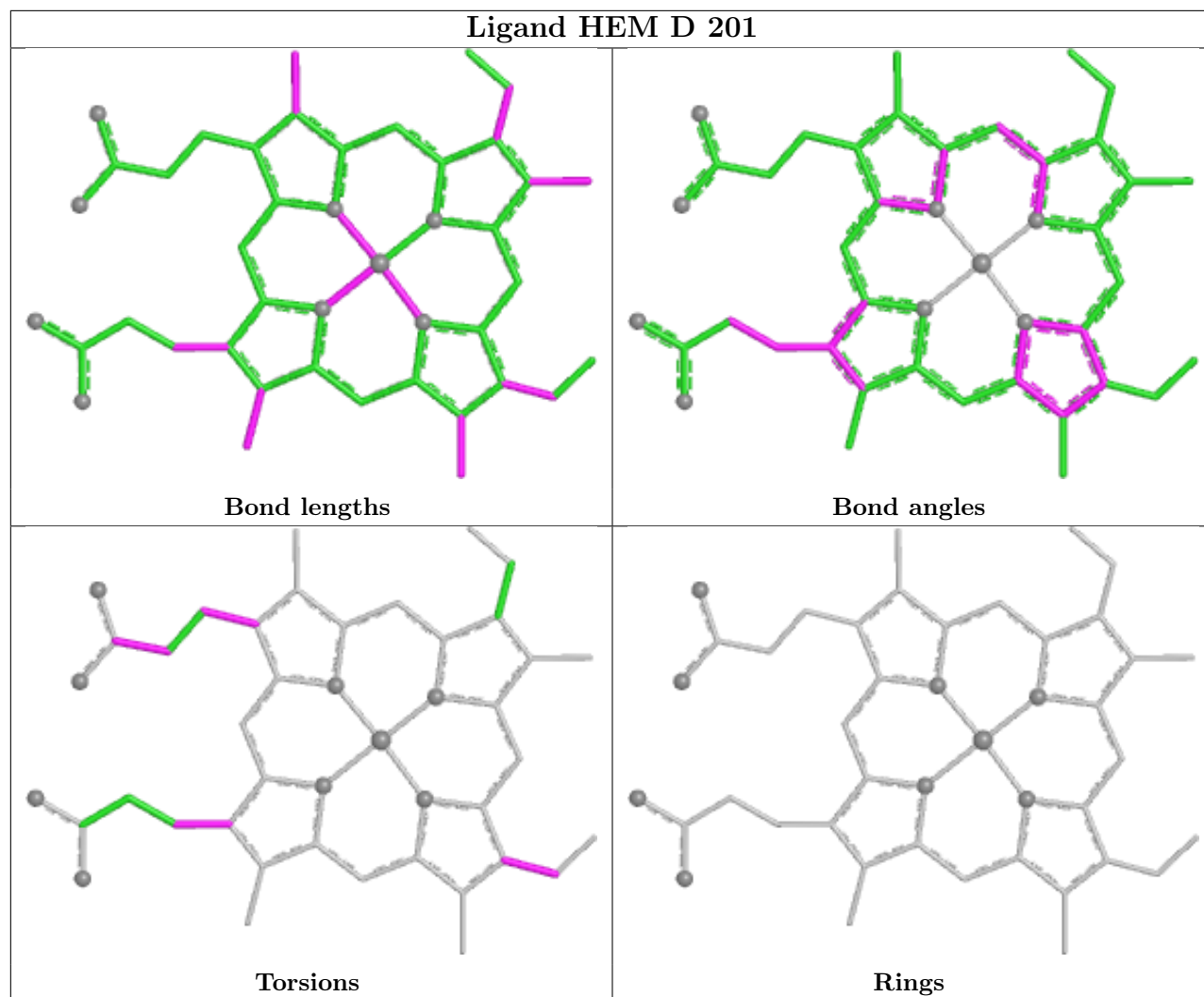


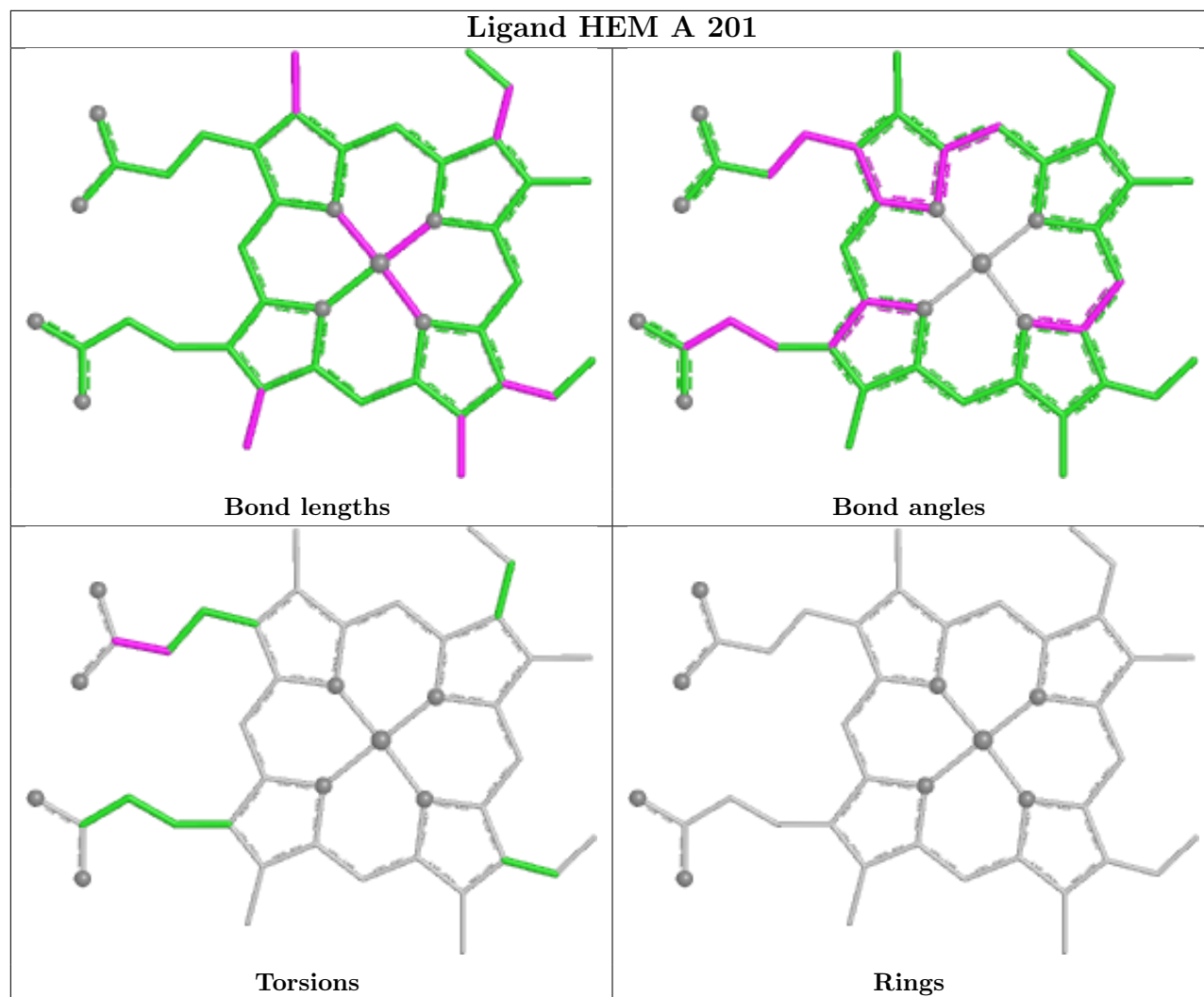
Ligand NBE D 203

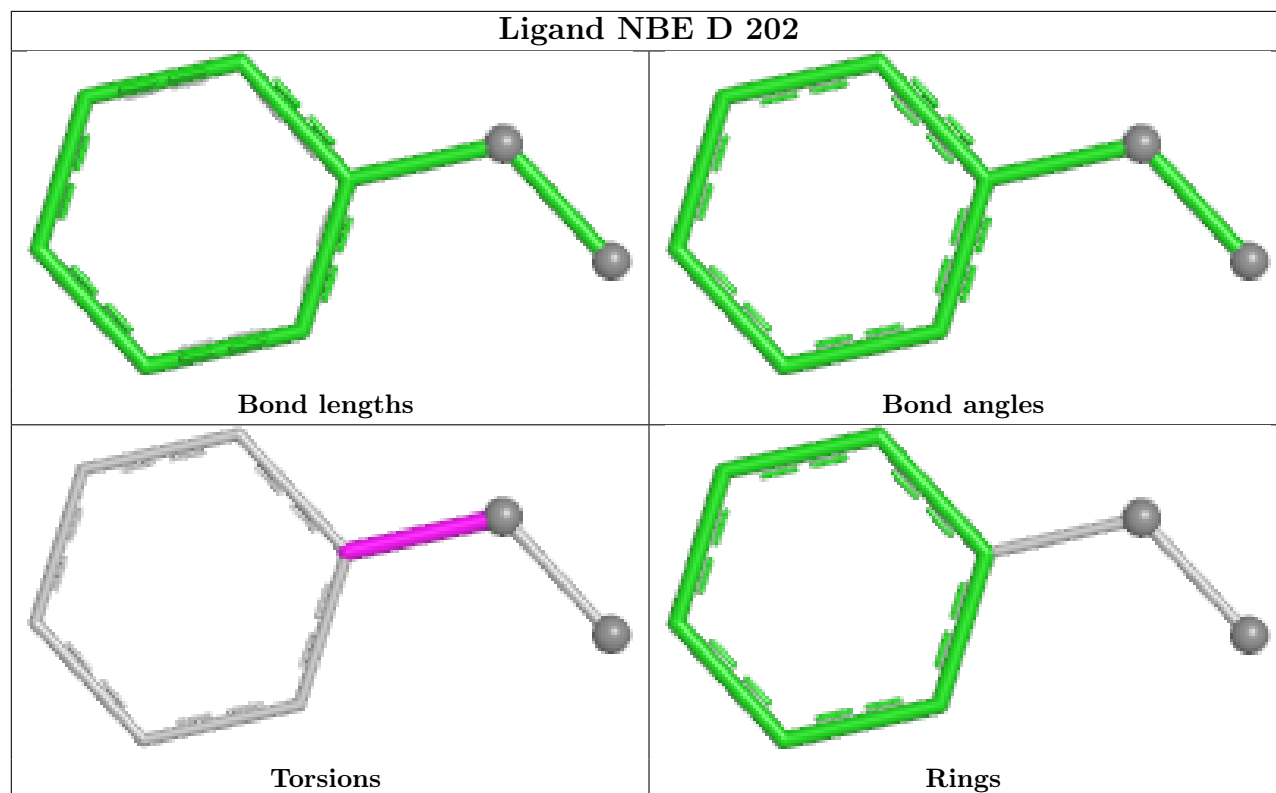


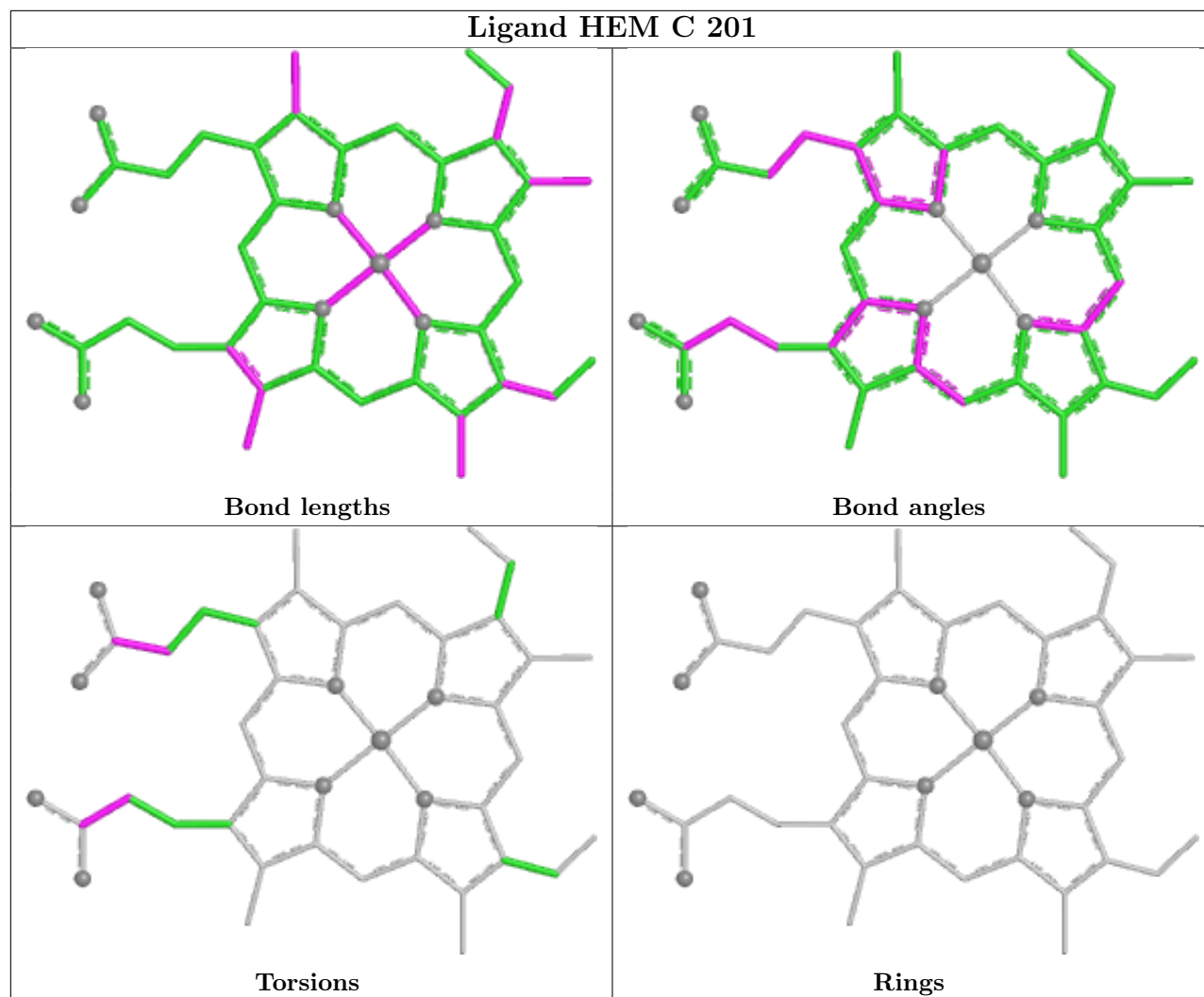
Ligand NBE C 202











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](#)

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	141/141 (100%)	-0.90	0 100 100	21, 30, 43, 63	0
1	C	141/141 (100%)	-0.88	0 100 100	20, 30, 43, 60	0
2	B	142/146 (97%)	-0.68	0 100 100	15, 38, 57, 71	3 (2%)
2	D	142/146 (97%)	-0.70	0 100 100	16, 39, 56, 68	2 (1%)
All	All	566/574 (98%)	-0.79	0 100 100	15, 34, 55, 71	5 (0%)

There are no RSRZ outliers to report.

6.2 Non-standard residues in protein, DNA, RNA chains [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	CSD	B	93	8/9	0.98	0.06	33,40,44,45	0
2	CSD	D	93	8/9	0.98	0.08	31,40,42,43	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

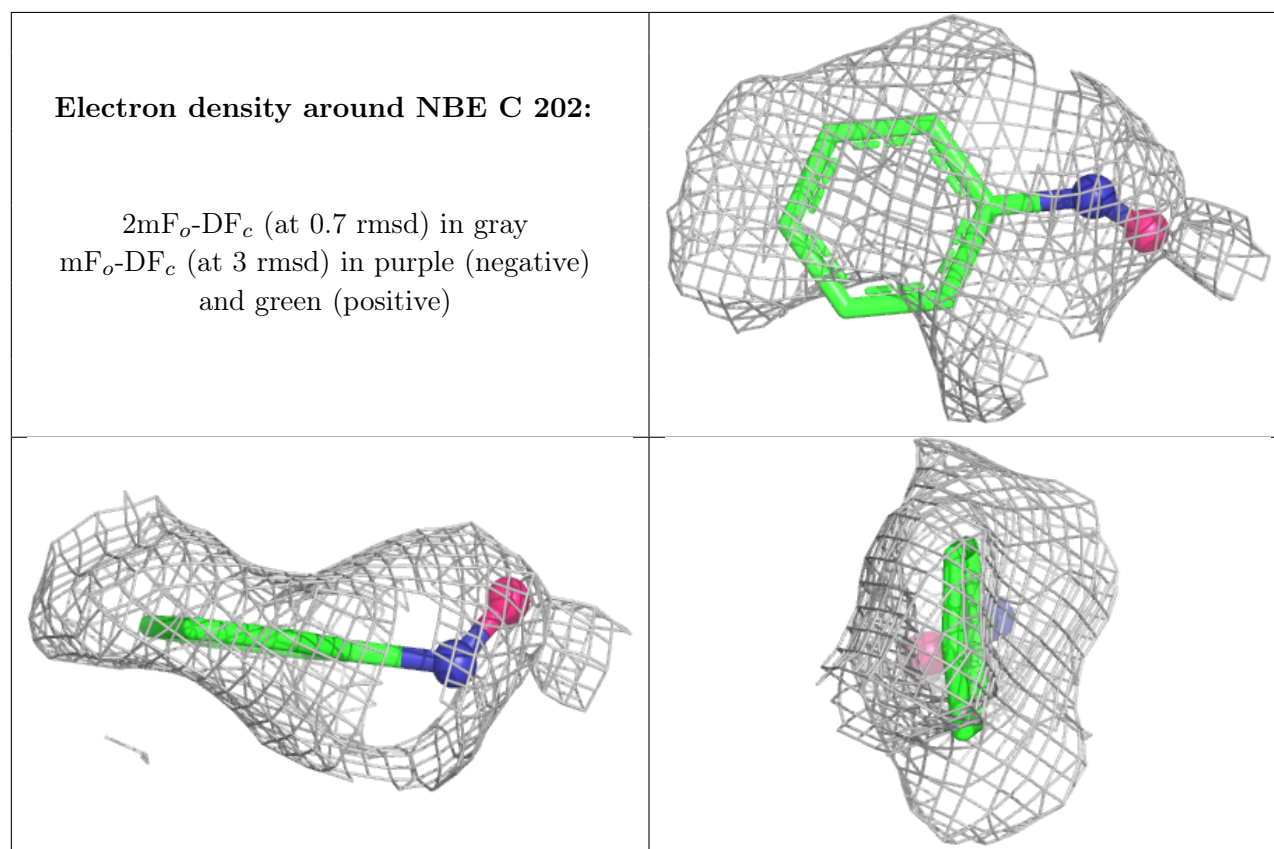
6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column

labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

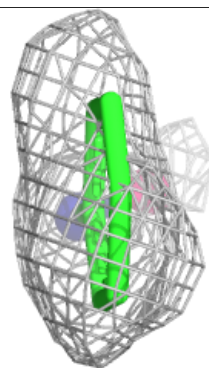
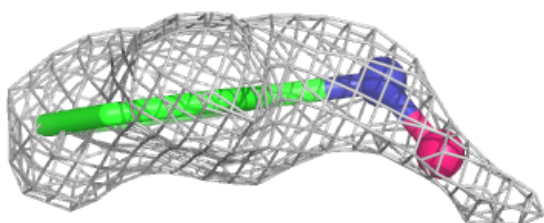
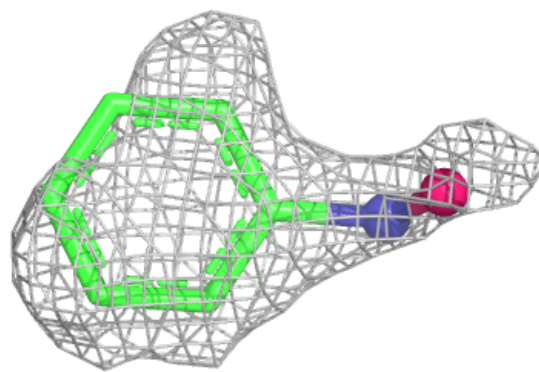
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	NBE	C	202	8/8	0.96	0.05	64,67,69,71	0
4	NBE	B	203	8/8	0.97	0.06	39,49,57,59	0
4	NBE	B	202	8/8	0.97	0.06	40,49,57,65	0
4	NBE	D	202	8/8	0.97	0.07	43,50,60,67	0
4	NBE	D	203	8/8	0.97	0.08	36,46,55,57	0
5	GOL	C	203	6/6	0.97	0.05	35,44,45,53	0
3	HEM	C	201	43/43	0.99	0.04	21,26,43,47	0
3	HEM	D	201	43/43	0.99	0.05	34,41,59,63	0
3	HEM	A	201	43/43	0.99	0.04	22,26,40,50	0
3	HEM	B	201	43/43	0.99	0.05	34,43,55,60	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.

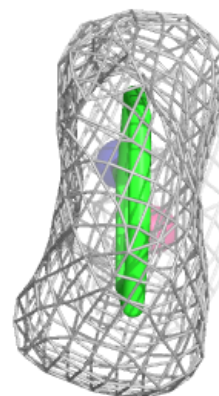
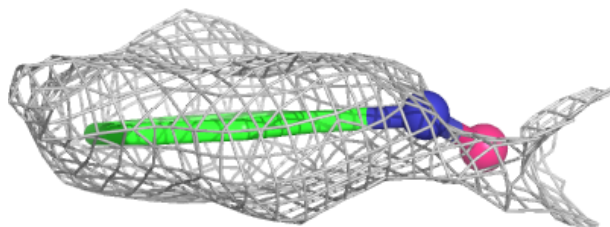
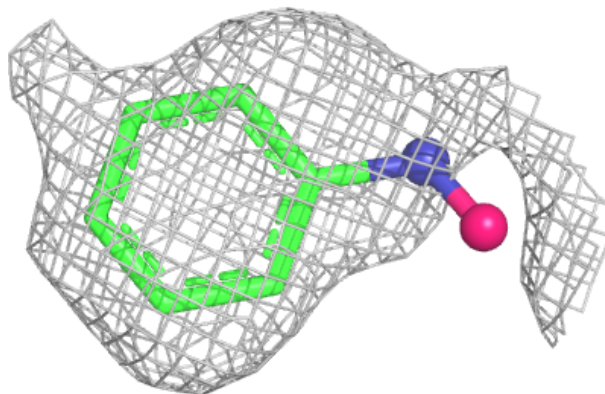


Electron density around NBE B 203:

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

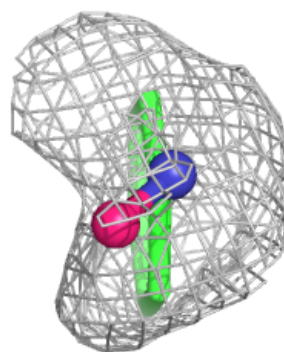
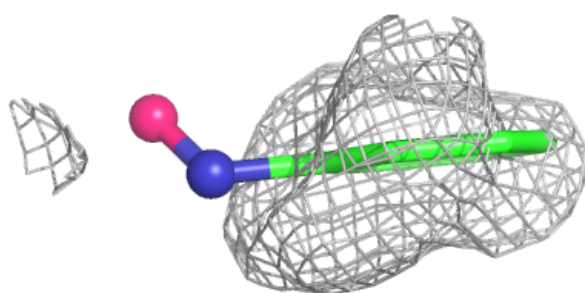
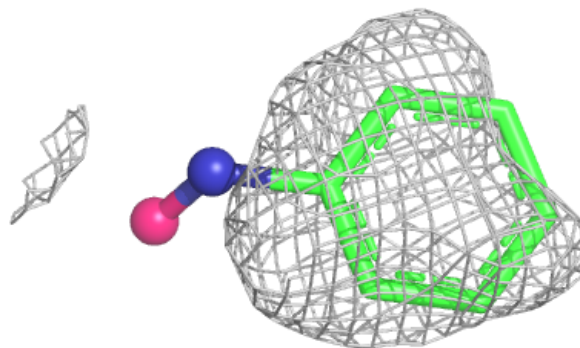
**Electron density around NBE B 202:**

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

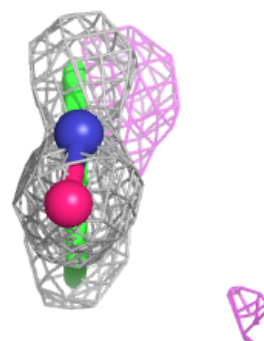
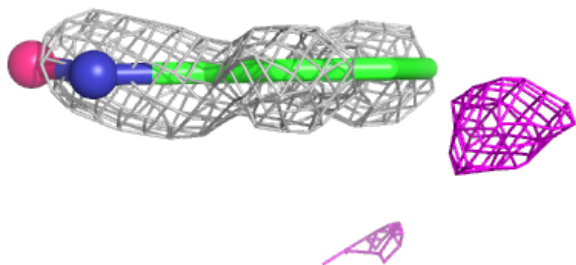
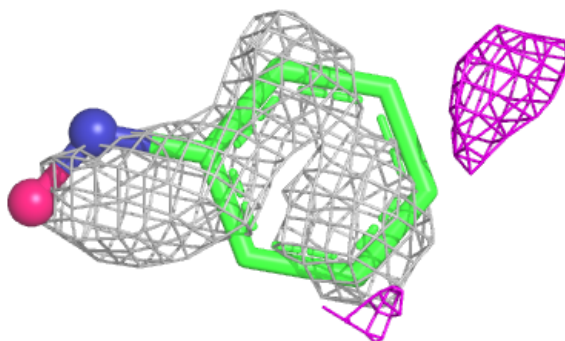


Electron density around NBE D 202:

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

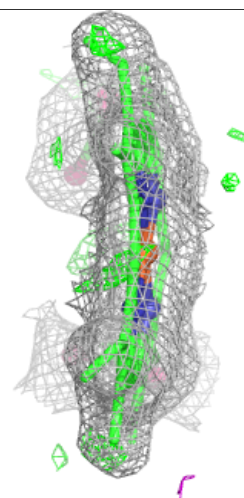
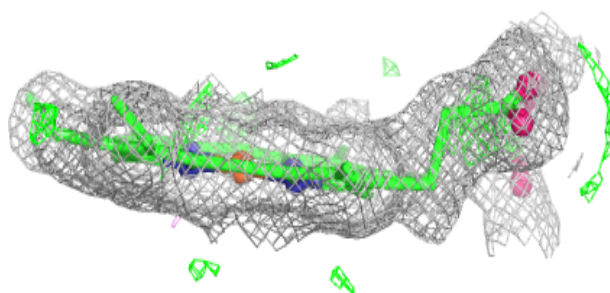
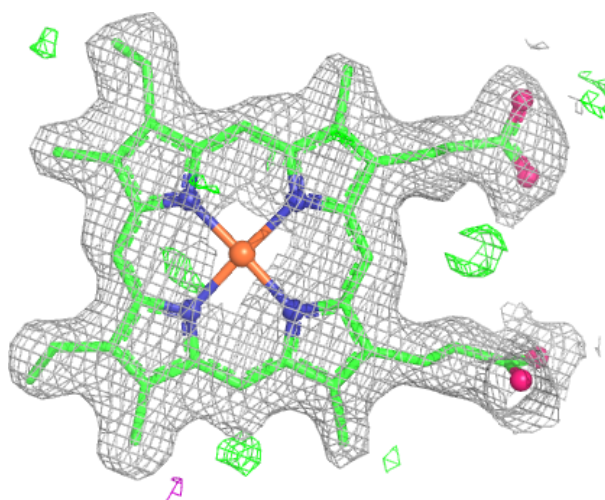
**Electron density around NBE D 203:**

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



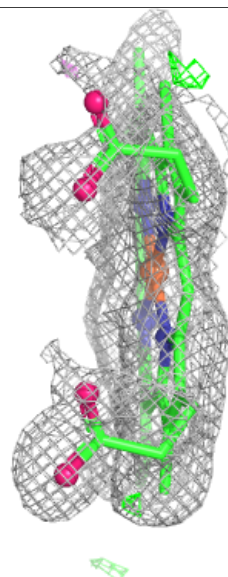
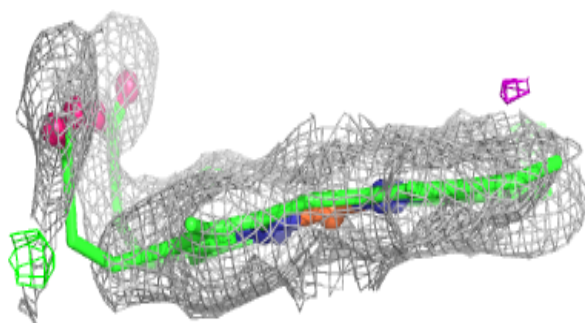
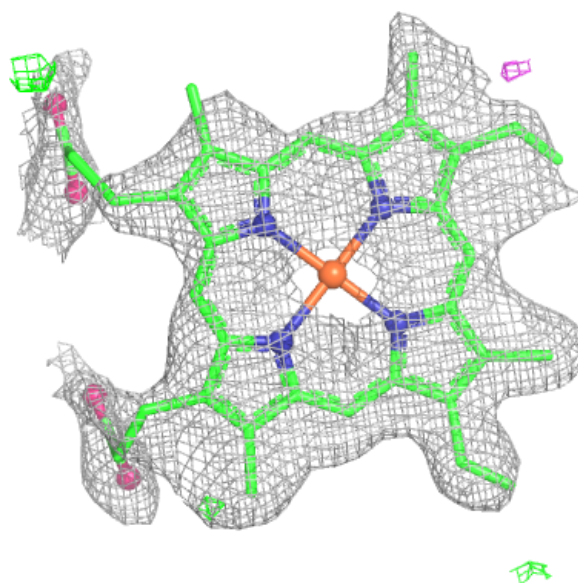
Electron density around HEM C 201:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



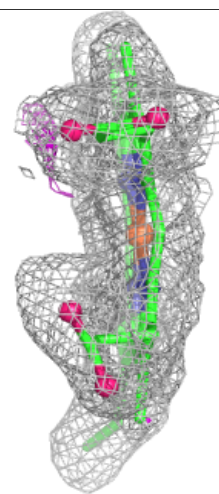
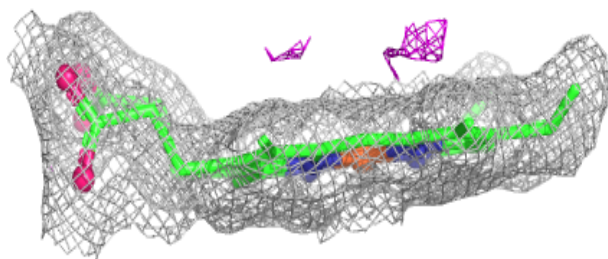
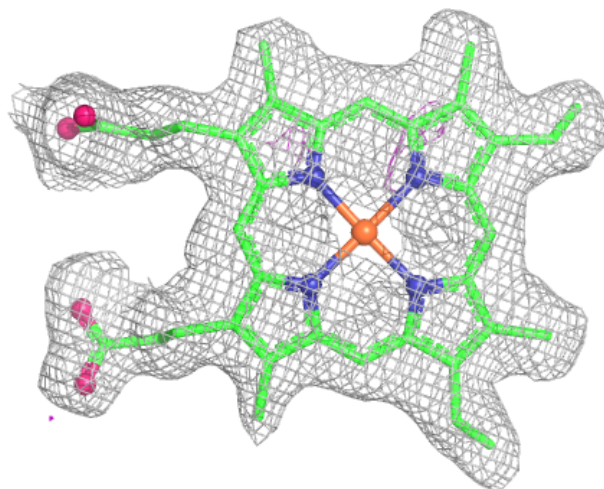
Electron density around HEM D 201:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



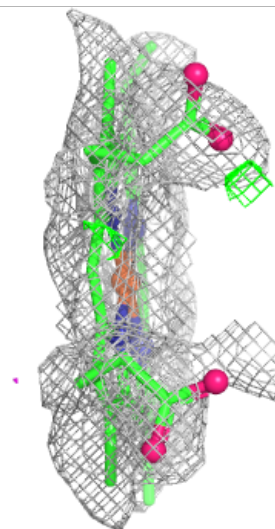
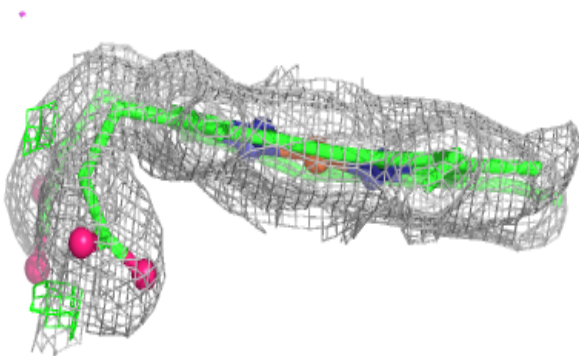
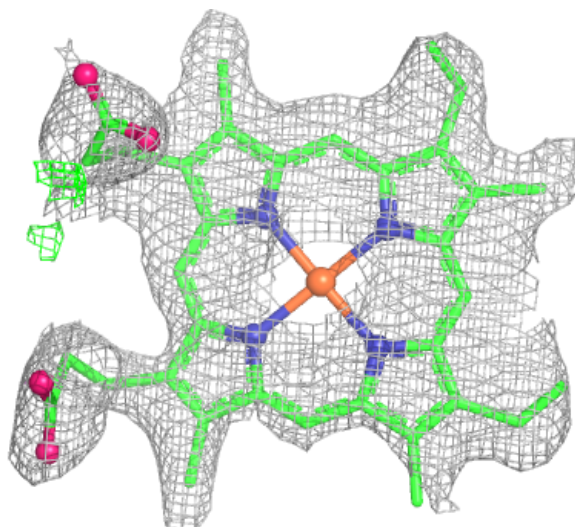
Electron density around HEM A 201:

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



Electron density around HEM B 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 [i](#)

There are no such residues in this entry.