



wwPDB X-ray Structure Validation Summary Report ⓘ

Sep 25, 2023 – 02:24 AM EDT

PDB ID : 5VVB
Title : Structure of human endothelial nitric oxide synthase heme domain in complex with 4-(2-(((2-Aminoquinolin-7-yl)methyl)amino)ethyl)-2-methylbenzotrile
Authors : Li, H.; Poulos, T.L.
Deposited on : 2017-05-19
Resolution : 2.15 Å(reported)

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

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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 : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtrriage (Phenix) : 1.13
EDS : 2.35.1
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) : 2.35.1

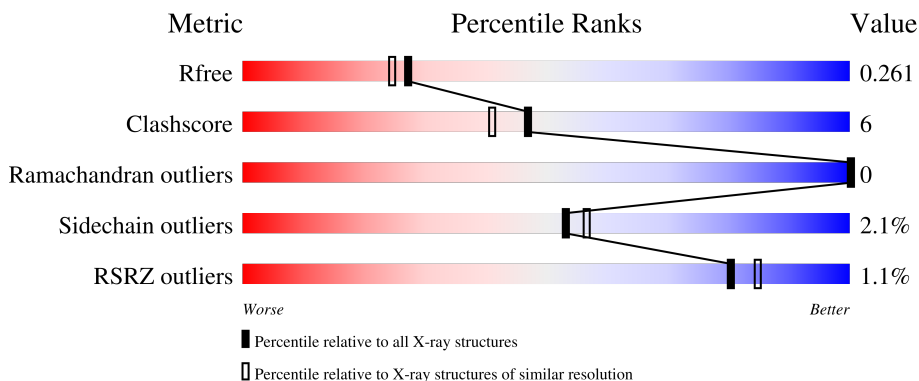
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.15 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	1479 (2.16-2.16)
Clashscore	141614	1585 (2.16-2.16)
Ramachandran outliers	138981	1560 (2.16-2.16)
Sidechain outliers	138945	1559 (2.16-2.16)
RSRZ outliers	127900	1456 (2.16-2.16)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	440	 2% 74% 15% 9%
1	B	440	 84% 8% 9%
1	C	440	 % 79% 12% 9%
1	D	440	 % 79% 12% 8%

2 Entry composition [i](#)

There are 10 unique types of molecules in this entry. The entry contains 13891 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 Nitric oxide synthase, endothelial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	401	Total 3212	C 2046	N 563	O 587	S 16	0	3	0
1	B	402	Total 3232	C 2057	N 568	O 590	S 17	0	5	0
1	C	401	Total 3212	C 2046	N 563	O 587	S 16	0	3	0
1	D	403	Total 3229	C 2056	N 569	O 588	S 16	0	3	0

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



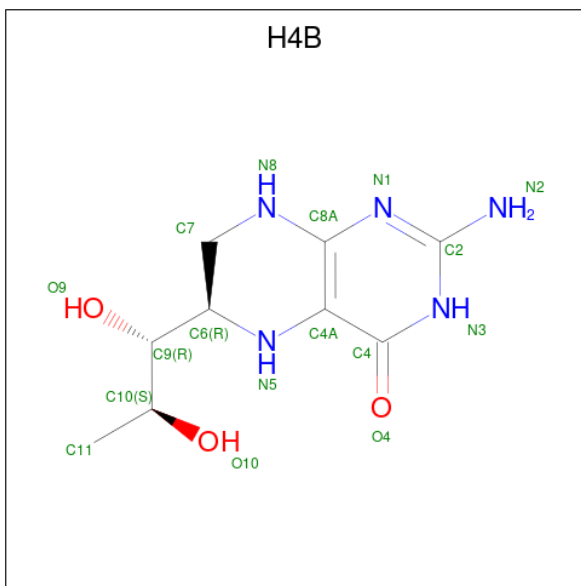
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
			Total	C	Fe	N			O
2	A	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
2	B	1	Total 43	C 34	Fe 1	N 4	O 4	0	0

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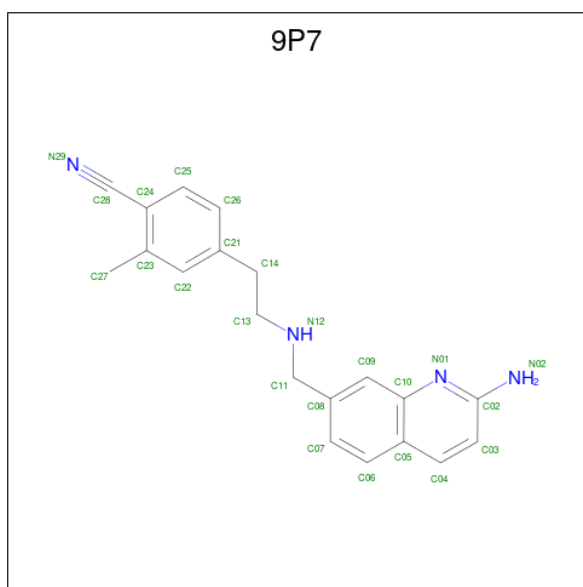
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
2	C	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
2	D	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		

- Molecule 3 is 5,6,7,8-TETRAHYDROBIOPTERIN (three-letter code: H4B) (formula: $C_9H_{15}N_5O_3$).



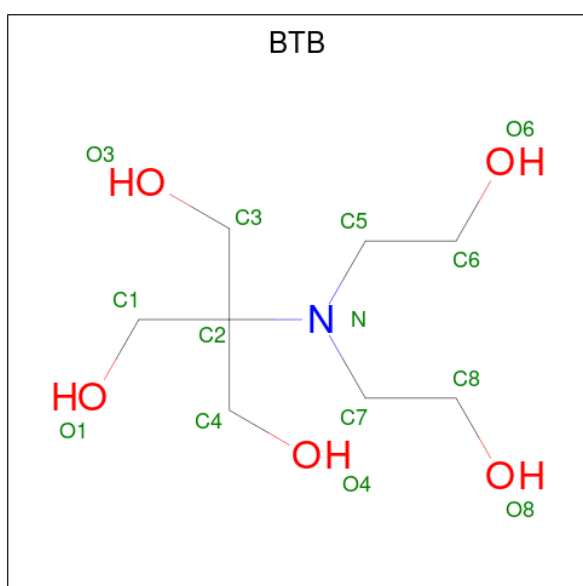
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			17	9	5	3		
3	B	1	Total	C	N	O	0	0
			17	9	5	3		
3	C	1	Total	C	N	O	0	0
			17	9	5	3		
3	D	1	Total	C	N	O	0	0
			17	9	5	3		

- Molecule 4 is 4-(2-[(2-aminoquinolin-7-yl)methyl]amino)ethyl)-2-methylbenzotrile (three-letter code: 9P7) (formula: $C_{20}H_{20}N_4$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	N	0	0
			24	20	4		
4	B	1	Total	C	N	0	0
			24	20	4		
4	C	1	Total	C	N	0	0
			24	20	4		
4	D	1	Total	C	N	0	0
			24	20	4		

- Molecule 5 is 2-[BIS-(2-HYDROXY-ETHYL)-AMINO]-2-HYDROXYMETHYL-PROPAN E-1,3-DIOL (three-letter code: BTB) (formula: C₈H₁₉NO₅).

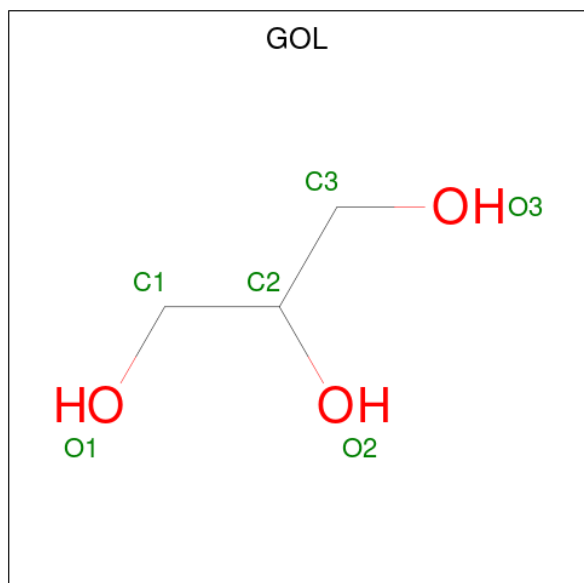


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	O	0	0
			14	8	1	5		
5	A	1	Total	C	N	O	0	0
			14	8	1	5		
5	B	1	Total	C	N	O	0	0
			14	8	1	5		
5	B	1	Total	C	N	O	0	0
			14	8	1	5		
5	B	1	Total	C	N	O	0	0
			14	8	1	5		
5	C	1	Total	C	N	O	0	0
			14	8	1	5		
5	C	1	Total	C	N	O	0	0
			14	8	1	5		
5	D	1	Total	C	N	O	0	0
			14	8	1	5		
5	D	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 6 is ZINC ION (three-letter code: ZN) (formula: Zn).

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

- Molecule 7 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



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

- Molecule 8 is GADOLINIUM ATOM (three-letter code: GD) (formula: Gd).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	1	Total	Gd	0	0
			1	1		
8	B	1	Total	Gd	0	0
			1	1		
8	C	1	Total	Gd	0	0
			1	1		
8	D	1	Total	Gd	0	0
			1	1		

- Molecule 9 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	1	Total	Cl	0	0
			1	1		
9	B	1	Total	Cl	0	0
			1	1		
9	C	1	Total	Cl	0	0
			1	1		
9	D	1	Total	Cl	0	0
			1	1		

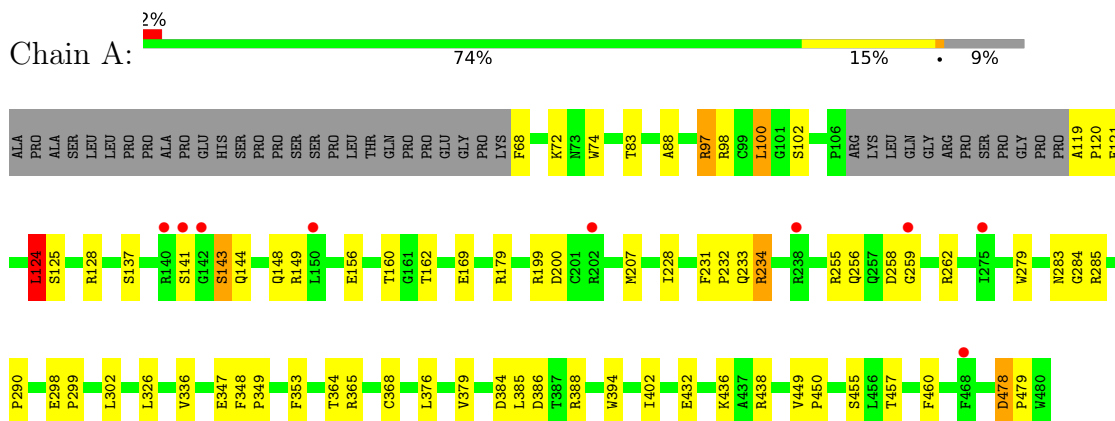
- Molecule 10 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	A	123	Total	O	0	0
			123	123		
10	B	148	Total	O	0	0
			148	148		
10	C	99	Total	O	0	0
			99	99		
10	D	152	Total	O	0	0
			152	152		

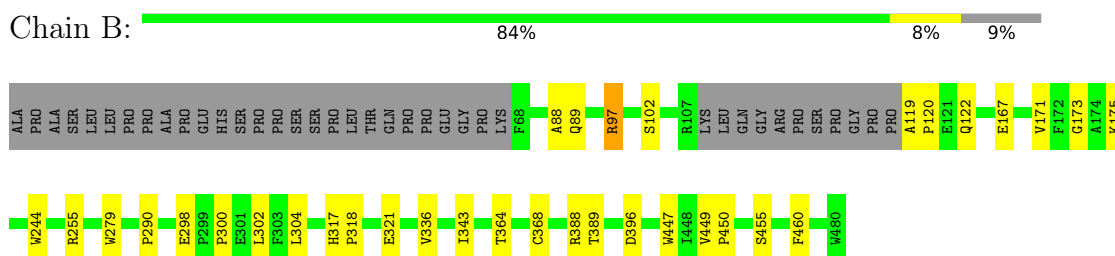
3 Residue-property plots [i](#)

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.

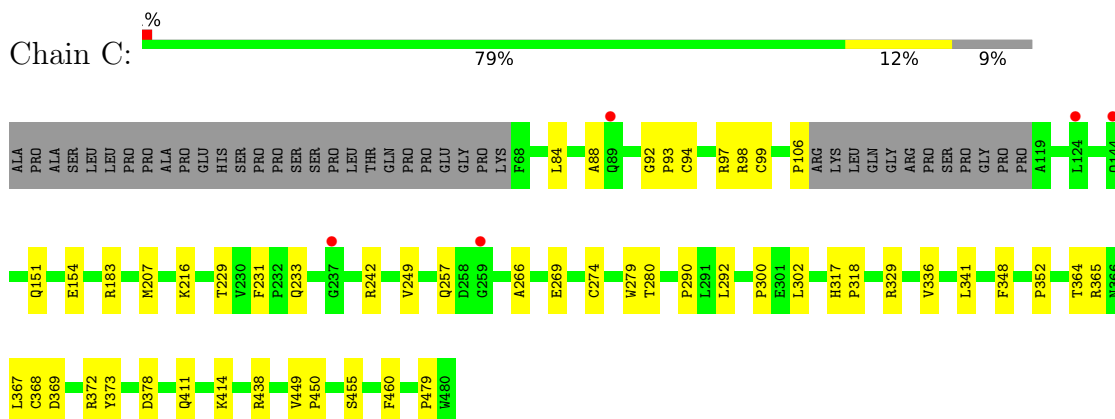
- Molecule 1: Nitric oxide synthase, endothelial




- Molecule 1: Nitric oxide synthase, endothelial

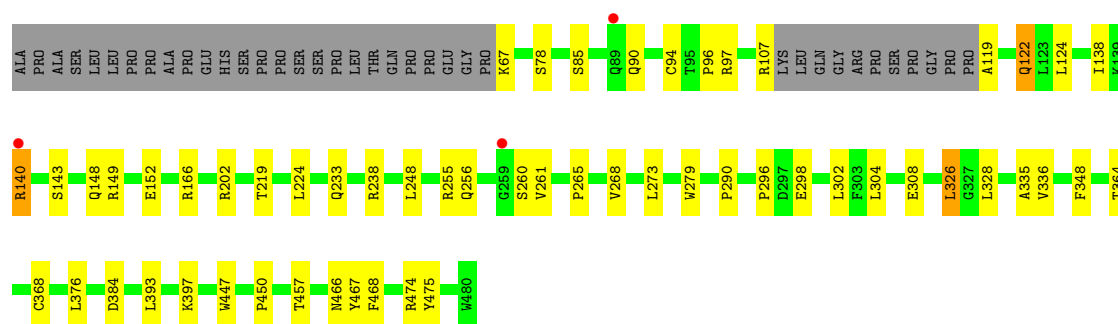


- Molecule 1: Nitric oxide synthase, endothelial



- Molecule 1: Nitric oxide synthase, endothelial

Chain D: 



4 Data and refinement statistics i

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	59.31Å 153.05Å 108.68Å 90.00° 90.75° 90.00°	Depositor
Resolution (Å)	55.30 – 2.15 55.30 – 2.15	Depositor EDS
% Data completeness (in resolution range)	92.5 (55.30-2.15) 91.8 (55.30-2.15)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	0.15	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.25 (at 2.14Å)	Xtrriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.202 , 0.268 0.196 , 0.261	Depositor DCC
R_{free} test set	4831 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	30.7	Xtrriage
Anisotropy	0.845	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 39.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	0.117 for h,-k,-l	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	13891	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.92% 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: 9P7, GD, CL, GOL, BTB, ZN, HEM, H4B

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	0/3313	0.52	1/4515 (0.0%)
1	B	0.43	0/3336	0.54	0/4546
1	C	0.36	0/3313	0.51	0/4515
1	D	0.41	0/3330	0.54	0/4536
All	All	0.39	0/13292	0.53	1/18112 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	124	LEU	CA-CB-CG	5.25	127.37	115.30

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	3212	0	3114	49	0
1	B	3232	0	3137	19	0
1	C	3212	0	3114	31	0
1	D	3229	0	3139	44	0
2	A	43	0	30	3	0
2	B	43	0	30	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	C	43	0	30	3	0
2	D	43	0	30	2	0
3	A	17	0	15	2	0
3	B	17	0	15	1	0
3	C	17	0	15	1	0
3	D	17	0	15	1	0
4	A	24	0	0	2	0
4	B	24	0	0	1	0
4	C	24	0	0	1	0
4	D	24	0	0	1	0
5	A	28	0	37	7	0
5	B	42	0	56	12	0
5	C	28	0	38	1	0
5	D	28	0	37	7	0
6	A	1	0	0	0	0
6	C	1	0	0	0	0
7	A	6	0	8	1	0
7	C	6	0	8	0	0
8	A	1	0	0	0	0
8	B	1	0	0	0	0
8	C	1	0	0	0	0
8	D	1	0	0	0	0
9	A	1	0	0	0	0
9	B	1	0	0	0	0
9	C	1	0	0	0	0
9	D	1	0	0	0	0
10	A	123	0	0	6	0
10	B	148	0	0	1	0
10	C	99	0	0	1	0
10	D	152	0	0	2	0
All	All	13891	0	12868	162	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 6.

The worst 5 of 162 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)
2:A:501:HEM:HBB2	2:A:501:HEM:HHC	1.69	0.74
1:B:279:TRP:HB2	1:B:302:LEU:HD21	1.73	0.71
5:D:504:BTB:H82	5:D:504:BTB:O6	1.91	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:298:GLU:HG3	1:A:299:PRO:HD2	1.74	0.70
1:A:347:GLU:OE2	10:A:601:HOH:O	2.10	0.69

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	400/440 (91%)	389 (97%)	11 (3%)	0	100	100
1	B	403/440 (92%)	391 (97%)	12 (3%)	0	100	100
1	C	400/440 (91%)	388 (97%)	12 (3%)	0	100	100
1	D	402/440 (91%)	386 (96%)	16 (4%)	0	100	100
All	All	1605/1760 (91%)	1554 (97%)	51 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	343/373 (92%)	334 (97%)	9 (3%)	46	47
1	B	346/373 (93%)	339 (98%)	7 (2%)	55	59
1	C	343/373 (92%)	338 (98%)	5 (2%)	65	69

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	D	345/373 (92%)	338 (98%)	7 (2%)	55	59
All	All	1377/1492 (92%)	1349 (98%)	28 (2%)	53	59

5 of 28 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	389	THR
1	D	326	LEU
1	C	249	VAL
1	D	140	ARG
1	C	216	LYS

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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 33 ligands modelled in this entry, 10 are monoatomic - leaving 23 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
4	9P7	B	503	-	26,26,26	2.24	2 (7%)	33,35,35	1.21	2 (6%)
5	BTB	B	506	-	13,13,13	0.45	0	7,16,16	0.56	0
3	H4B	C	502	-	16,18,18	0.80	0	11,26,26	2.76	6 (54%)
3	H4B	A	502	-	16,18,18	1.07	0	11,26,26	2.70	5 (45%)
5	BTB	A	505	-	13,13,13	0.66	0	7,16,16	1.36	0
4	9P7	A	503	-	26,26,26	2.22	1 (3%)	33,35,35	1.32	4 (12%)
5	BTB	C	505	-	13,13,13	0.36	0	7,16,16	0.90	0
3	H4B	D	502	-	16,18,18	0.97	0	11,26,26	2.74	6 (54%)
7	GOL	C	507	-	5,5,5	0.38	0	5,5,5	0.30	0
5	BTB	D	504	8	13,13,13	0.41	0	7,16,16	0.76	0
2	HEM	B	501	1	41,50,50	1.95	6 (14%)	45,82,82	2.06	13 (28%)
5	BTB	D	505	-	13,13,13	0.49	0	7,16,16	0.72	0
3	H4B	B	502	-	16,18,18	0.85	0	11,26,26	2.68	6 (54%)
5	BTB	A	504	8	13,13,13	0.36	0	7,16,16	0.55	0
5	BTB	B	505	-	13,13,13	0.37	0	7,16,16	0.47	0
4	9P7	C	503	-	26,26,26	2.24	1 (3%)	33,35,35	1.29	7 (21%)
4	9P7	D	503	-	26,26,26	2.25	2 (7%)	33,35,35	1.27	4 (12%)
5	BTB	C	504	8	13,13,13	0.40	0	7,16,16	0.59	0
7	GOL	A	507	-	5,5,5	0.32	0	5,5,5	0.32	0
2	HEM	A	501	1	41,50,50	2.03	5 (12%)	45,82,82	1.88	8 (17%)
5	BTB	B	504	8	13,13,13	0.40	0	7,16,16	0.99	1 (14%)
2	HEM	C	501	1	41,50,50	1.95	8 (19%)	45,82,82	1.72	11 (24%)
2	HEM	D	501	1	41,50,50	1.94	6 (14%)	45,82,82	1.99	12 (26%)

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
4	9P7	B	503	-	-	0/9/18/18	0/3/3/3
5	BTB	B	506	-	-	14/21/21/21	-
3	H4B	C	502	-	-	3/8/17/17	0/2/2/2
3	H4B	A	502	-	-	3/8/17/17	0/2/2/2
5	BTB	A	505	-	-	8/21/21/21	-
4	9P7	A	503	-	-	2/9/18/18	0/3/3/3
5	BTB	C	505	-	-	8/21/21/21	-
3	H4B	D	502	-	-	3/8/17/17	0/2/2/2

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
7	GOL	C	507	-	-	3/4/4/4	-
5	BTB	D	504	8	-	4/21/21/21	-
2	HEM	B	501	1	-	3/12/54/54	-
5	BTB	D	505	-	-	11/21/21/21	-
3	H4B	B	502	-	-	3/8/17/17	0/2/2/2
5	BTB	A	504	8	-	7/21/21/21	-
5	BTB	B	505	-	-	8/21/21/21	-
4	9P7	C	503	-	-	1/9/18/18	0/3/3/3
4	9P7	D	503	-	-	1/9/18/18	0/3/3/3
5	BTB	C	504	8	-	4/21/21/21	-
7	GOL	A	507	-	-	2/4/4/4	-
2	HEM	A	501	1	-	4/12/54/54	-
5	BTB	B	504	8	-	4/21/21/21	-
2	HEM	C	501	1	-	5/12/54/54	-
2	HEM	D	501	1	-	0/12/54/54	-

The worst 5 of 31 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	D	503	9P7	C24-C28	-10.36	1.28	1.44
4	A	503	9P7	C24-C28	-10.32	1.28	1.44
4	C	503	9P7	C24-C28	-10.29	1.28	1.44
4	B	503	9P7	C24-C28	-10.25	1.28	1.44
2	A	501	HEM	C3D-C2D	7.98	1.53	1.36

The worst 5 of 85 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	501	HEM	C4D-ND-C1D	6.49	111.78	105.07
2	B	501	HEM	C4D-ND-C1D	6.21	111.49	105.07
3	A	502	H4B	C8A-C4A-C4	6.12	120.01	114.57
3	C	502	H4B	C8A-C4A-C4	5.84	119.76	114.57
2	B	501	HEM	C4C-CHD-C1D	5.83	130.25	122.56

There are no chirality outliers.

5 of 101 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	502	H4B	C7-C6-C9-O9

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Mol	Chain	Res	Type	Atoms
3	A	502	H4B	C7-C6-C9-C10
3	B	502	H4B	C7-C6-C9-O9
3	B	502	H4B	C7-C6-C9-C10
3	C	502	H4B	C7-C6-C9-O9

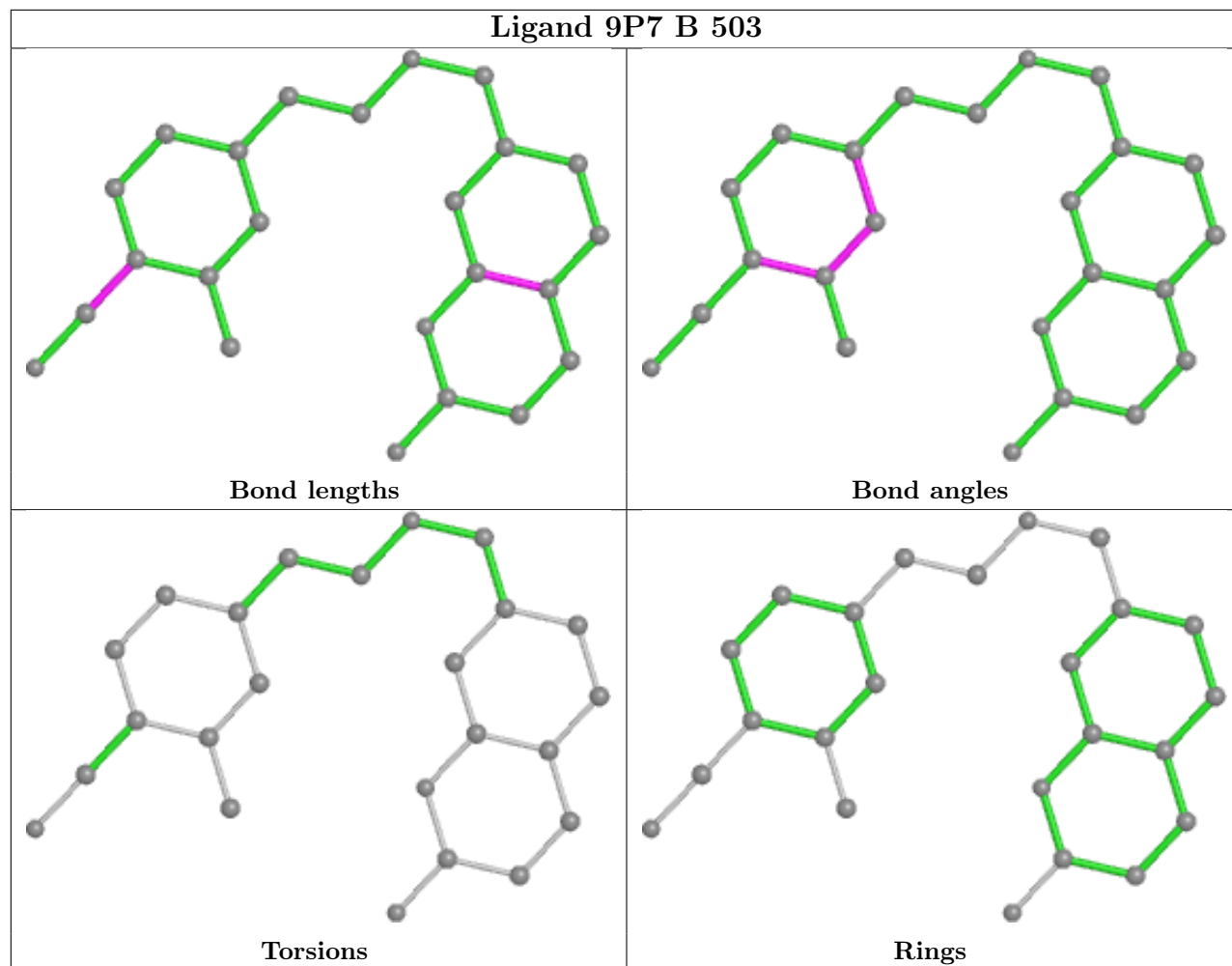
There are no ring outliers.

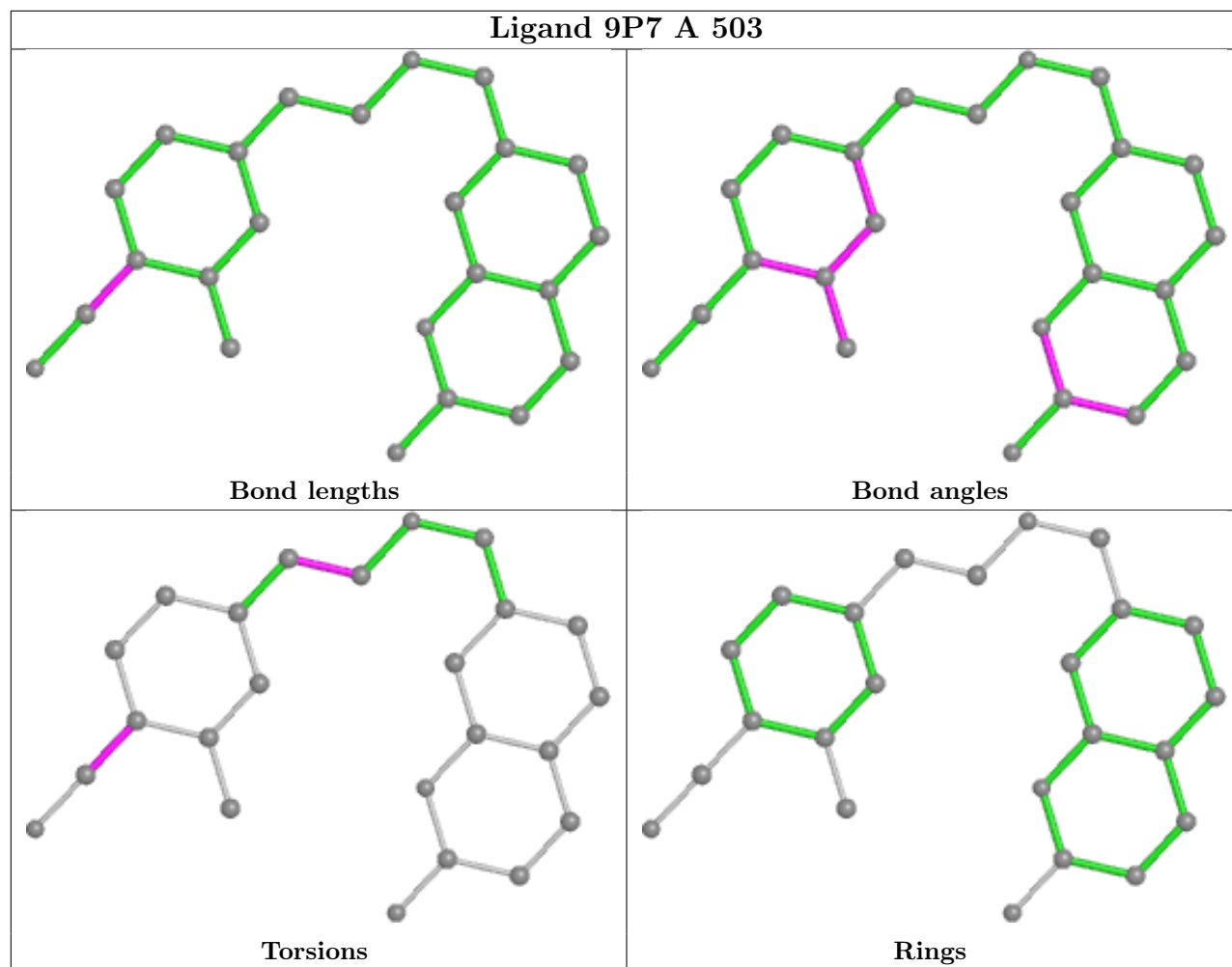
21 monomers are involved in 46 short contacts:

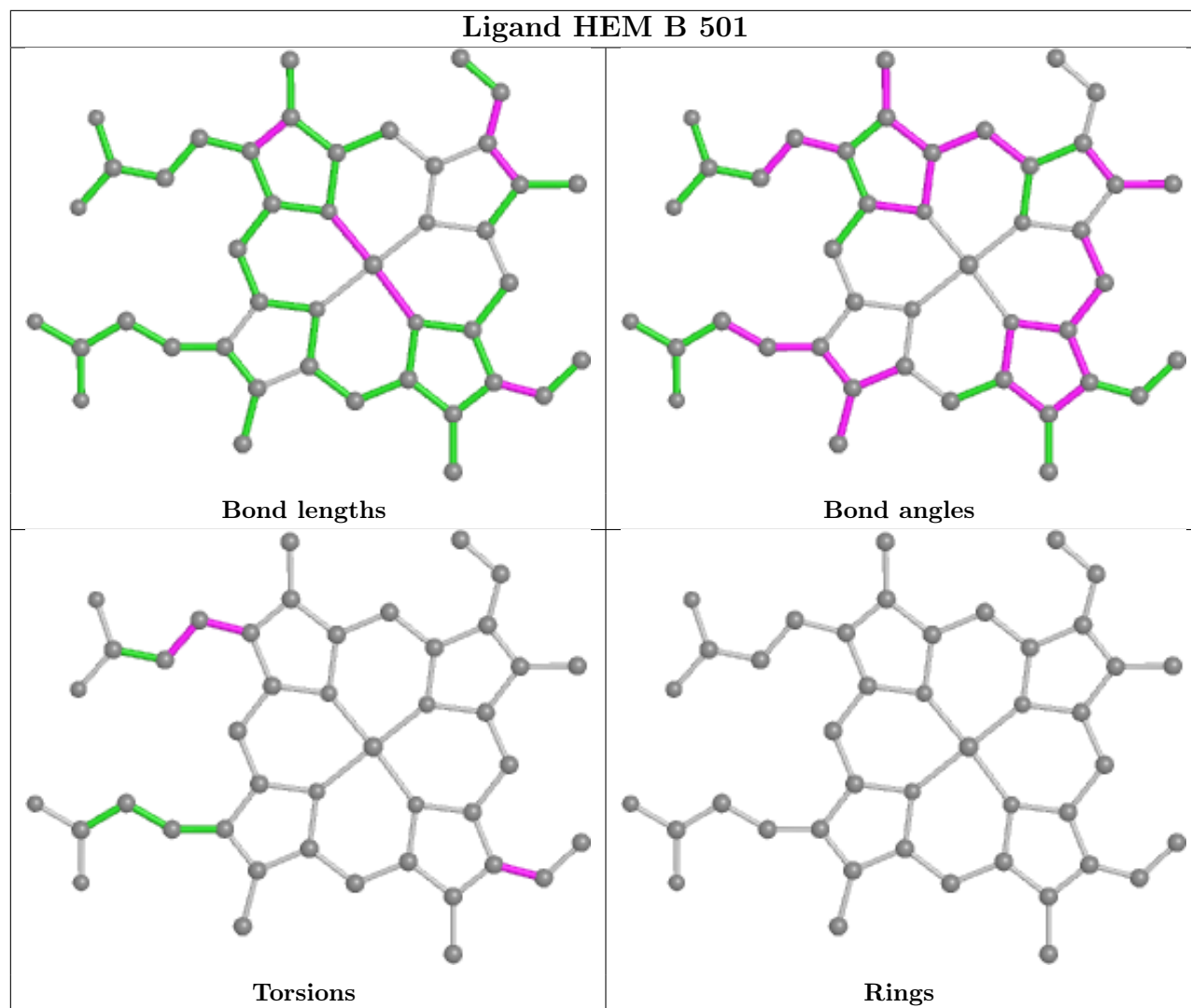
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	503	9P7	1	0
5	B	506	BTB	5	0
3	C	502	H4B	1	0
3	A	502	H4B	2	0
5	A	505	BTB	3	0
4	A	503	9P7	2	0
5	C	505	BTB	1	0
3	D	502	H4B	1	0
5	D	504	BTB	3	0
2	B	501	HEM	1	0
5	D	505	BTB	4	0
3	B	502	H4B	1	0
5	A	504	BTB	4	0
5	B	505	BTB	4	0
4	C	503	9P7	1	0
4	D	503	9P7	1	0
7	A	507	GOL	1	0
2	A	501	HEM	3	0
5	B	504	BTB	3	0
2	C	501	HEM	3	0
2	D	501	HEM	2	0

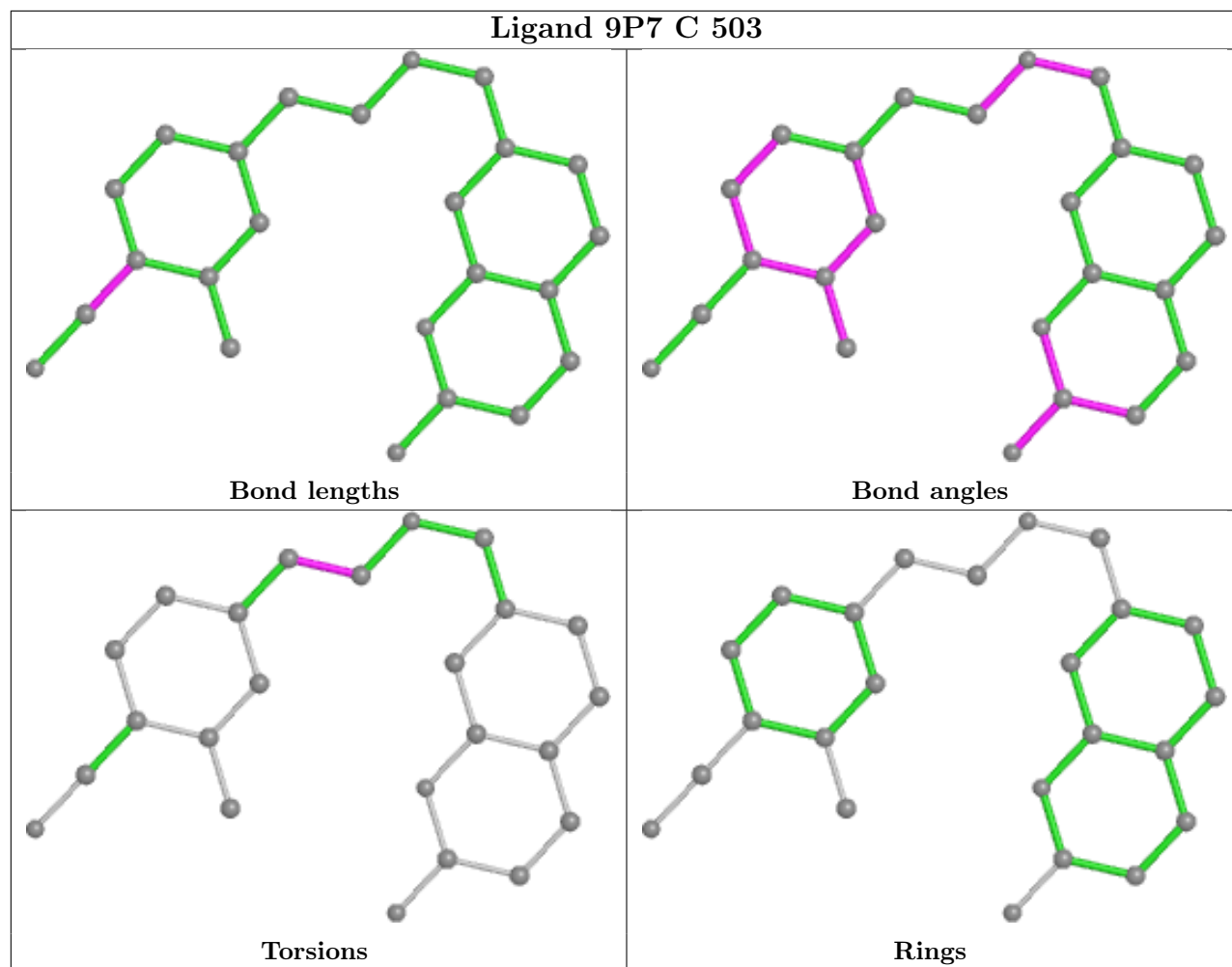
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient

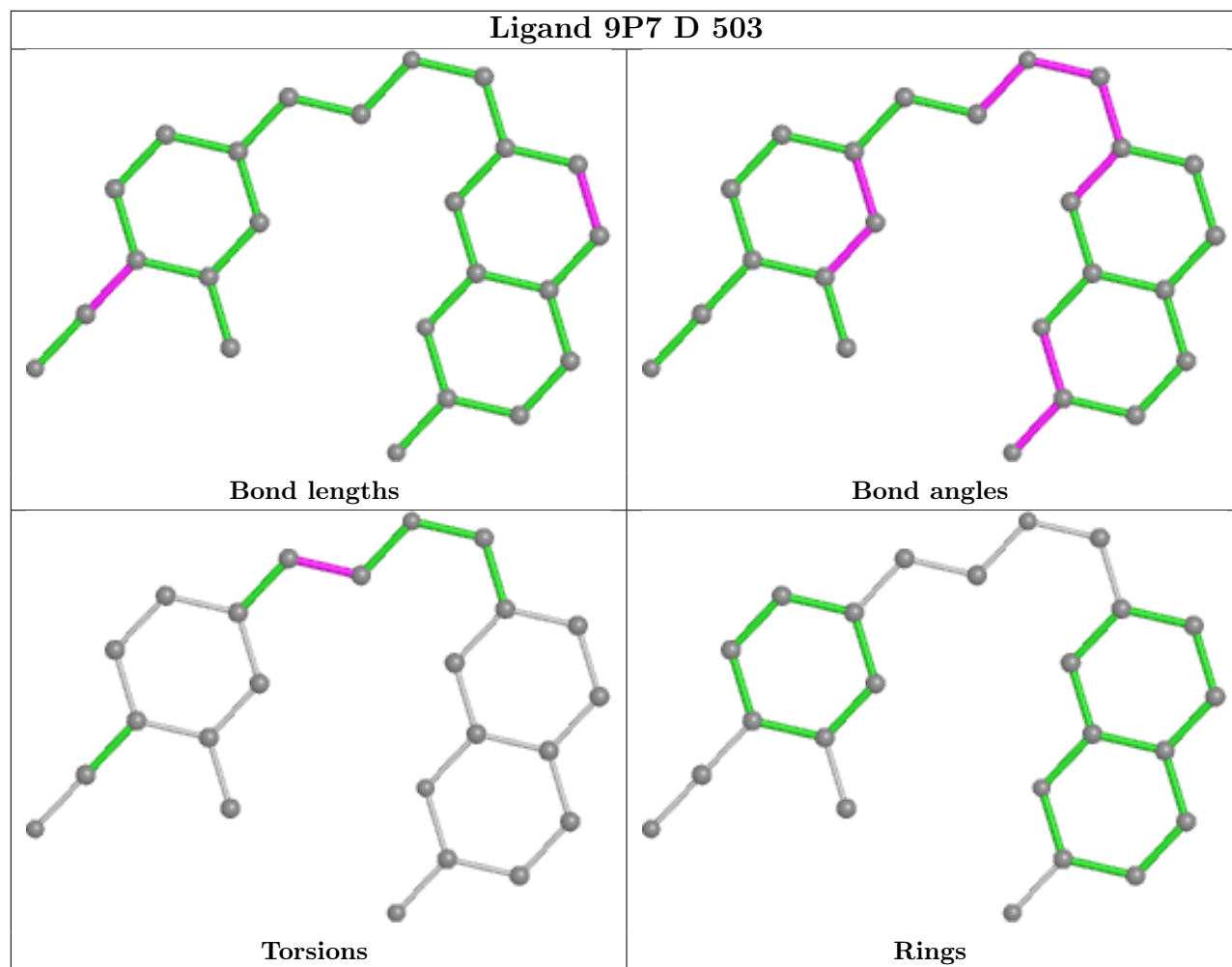
equivalents in the CSD to analyse the geometry.

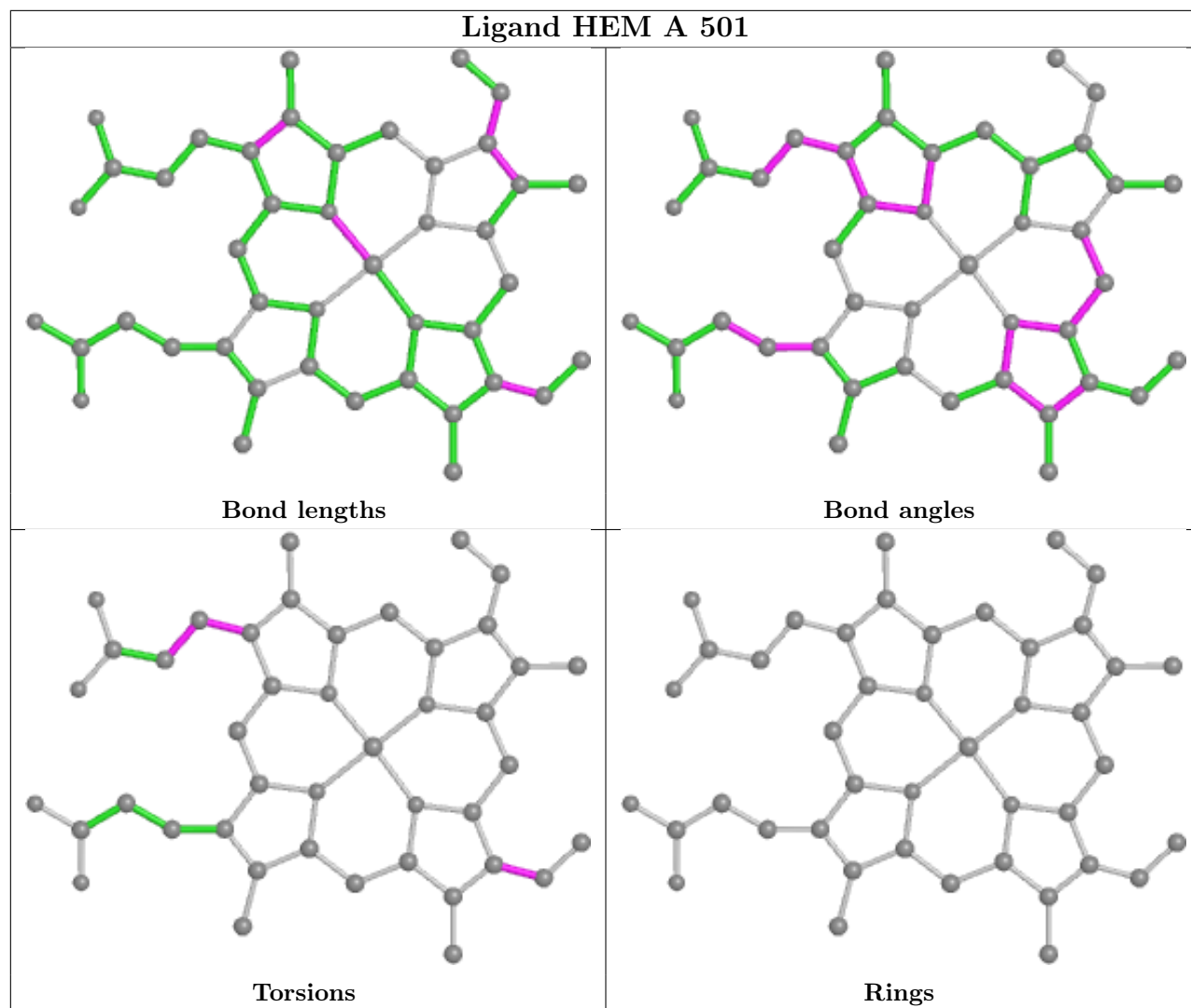


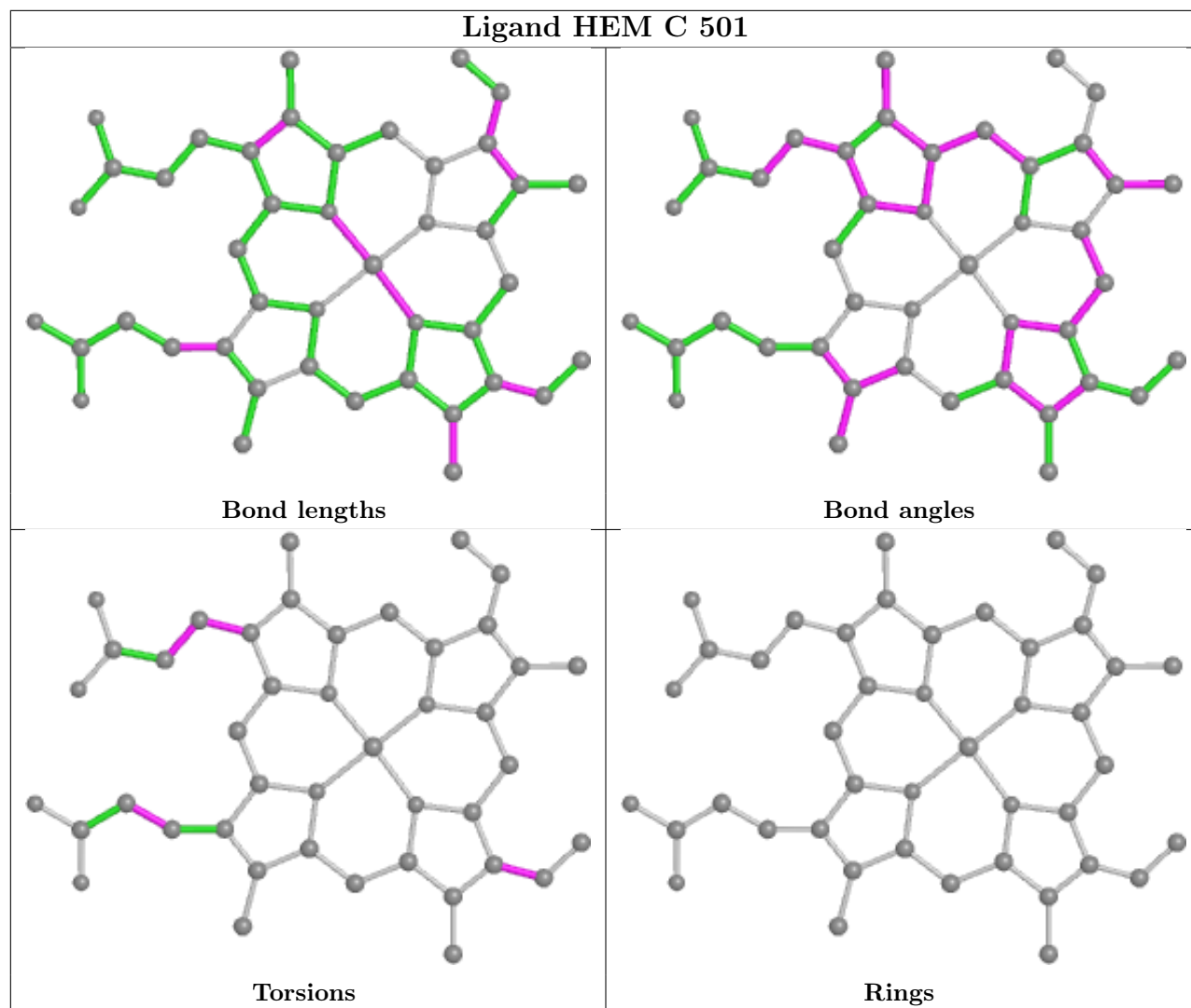


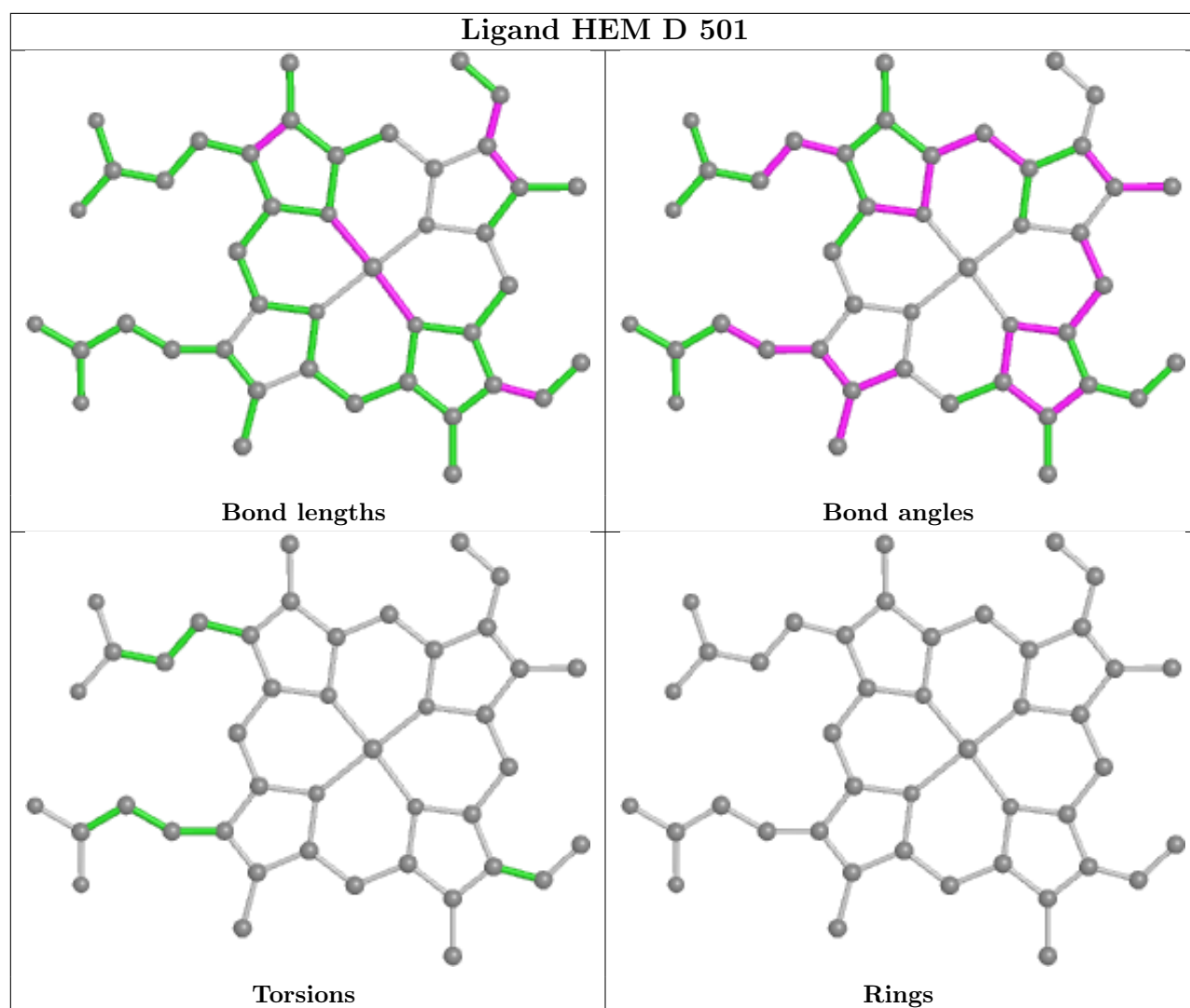












5.7 Other polymers [i](#)

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	401/440 (91%)	-0.23	9 (2%) 62 69	25, 47, 93, 120	0
1	B	402/440 (91%)	-0.47	0 100 100	20, 35, 65, 114	0
1	C	401/440 (91%)	-0.21	5 (1%) 79 83	23, 49, 93, 129	0
1	D	403/440 (91%)	-0.44	3 (0%) 87 91	20, 36, 75, 132	0
All	All	1607/1760 (91%)	-0.34	17 (1%) 80 85	20, 41, 88, 132	0

The worst 5 of 17 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	238	ARG	4.6
1	C	89	GLN	3.4
1	A	140	ARG	3.4
1	A	468	PHE	3.2
1	D	89	GLN	3.1

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

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

6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

6.4 Ligands [i](#)

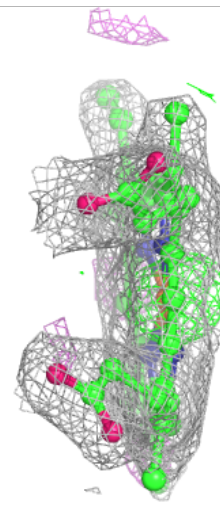
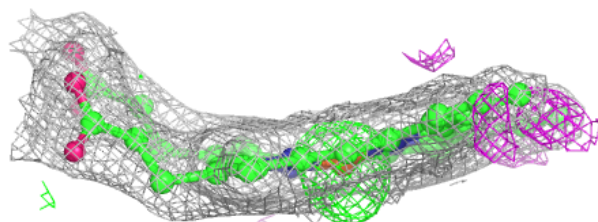
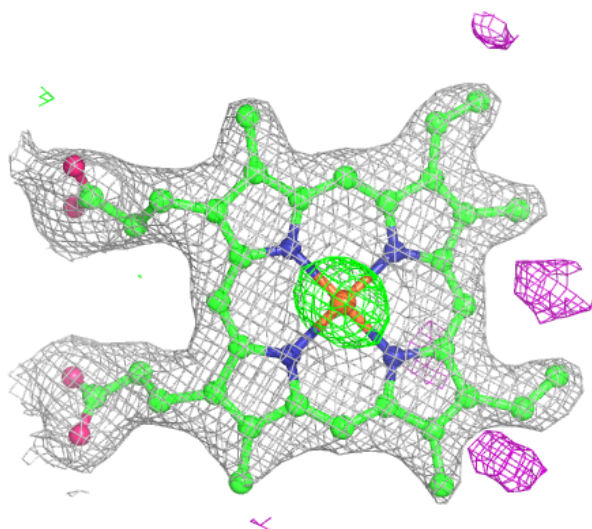
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q < 0.9’ lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	BTB	C	505	14/14	0.85	0.16	69,76,84,88	0
5	BTB	B	505	14/14	0.87	0.17	66,76,92,95	0
5	BTB	D	505	14/14	0.89	0.23	70,78,92,99	0
8	GD	C	508	1/1	0.89	0.11	131,131,131,131	0
7	GOL	C	507	6/6	0.90	0.14	38,51,77,82	0
5	BTB	B	504	14/14	0.91	0.11	30,60,66,70	0
5	BTB	B	506	14/14	0.91	0.08	60,67,76,78	0
5	BTB	A	505	14/14	0.92	0.18	58,71,76,84	0
5	BTB	C	504	14/14	0.92	0.17	70,84,95,100	0
5	BTB	D	504	14/14	0.93	0.11	15,44,59,66	0
7	GOL	A	507	6/6	0.93	0.16	42,51,66,66	0
9	CL	C	509	1/1	0.93	0.12	66,66,66,66	0
2	HEM	C	501	43/43	0.94	0.14	29,44,58,87	0
4	9P7	C	503	24/24	0.95	0.12	30,44,54,76	0
3	H4B	D	502	17/17	0.96	0.10	29,34,44,45	0
4	9P7	A	503	24/24	0.96	0.09	26,38,47,61	0
3	H4B	A	502	17/17	0.96	0.12	24,45,67,68	0
2	HEM	B	501	43/43	0.97	0.10	12,25,39,65	0
3	H4B	B	502	17/17	0.97	0.11	24,34,43,52	0
4	9P7	B	503	24/24	0.97	0.11	22,31,47,56	0
3	H4B	C	502	17/17	0.97	0.10	28,43,59,67	0
8	GD	A	508	1/1	0.97	0.11	92,92,92,92	0
4	9P7	D	503	24/24	0.97	0.08	17,30,40,42	0
5	BTB	A	504	14/14	0.97	0.13	15,75,84,96	0
2	HEM	A	501	43/43	0.99	0.10	24,38,47,59	0
9	CL	A	509	1/1	0.99	0.10	65,65,65,65	0
2	HEM	D	501	43/43	0.99	0.09	13,23,38,55	0
9	CL	D	507	1/1	0.99	0.07	37,37,37,37	0
8	GD	D	506	1/1	1.00	0.16	43,43,43,43	0
6	ZN	A	506	1/1	1.00	0.10	26,26,26,26	0
9	CL	B	508	1/1	1.00	0.09	34,34,34,34	0
8	GD	B	507	1/1	1.00	0.16	45,45,45,45	0
6	ZN	C	506	1/1	1.00	0.13	36,36,36,36	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.

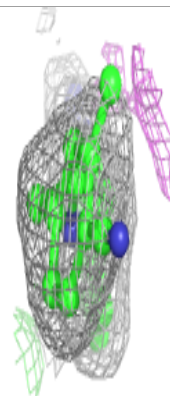
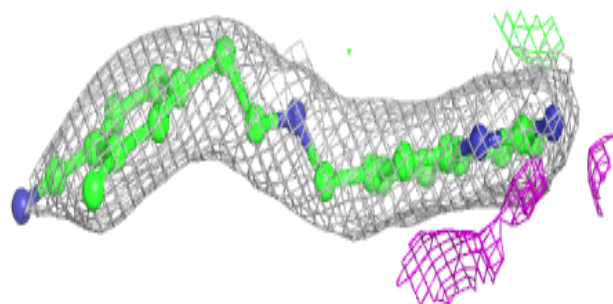
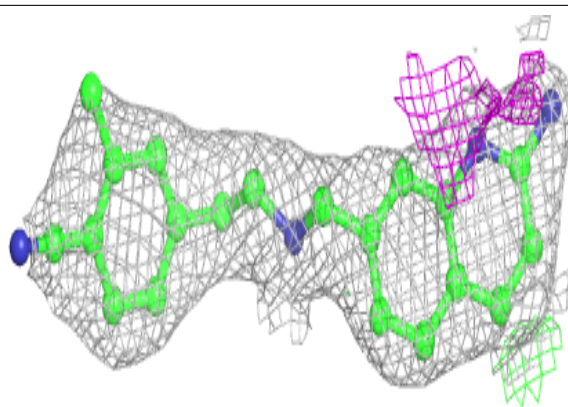
Electron density around HEM C 501:

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

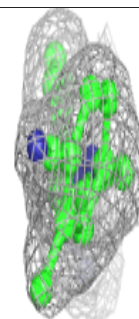
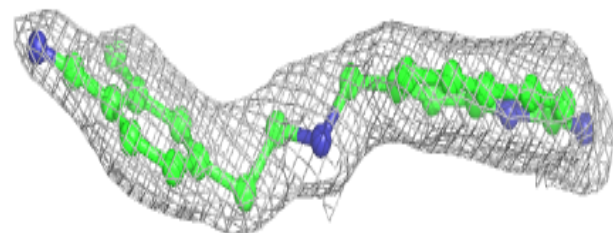
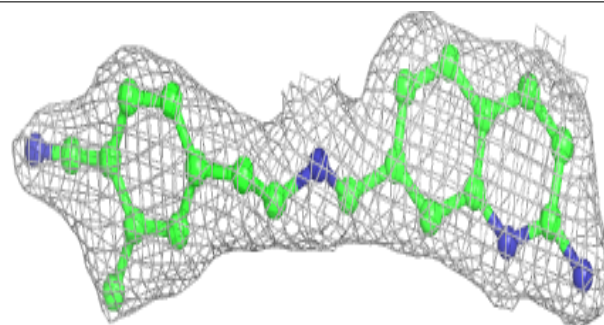


Electron density around 9P7 C 503:

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

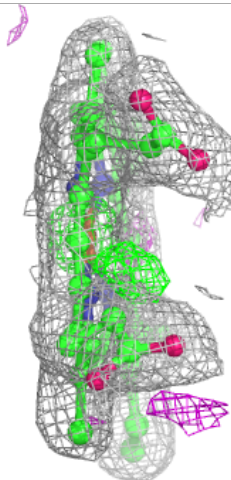
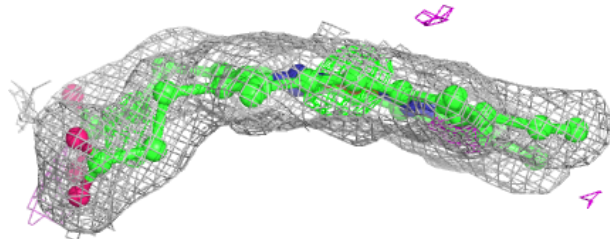
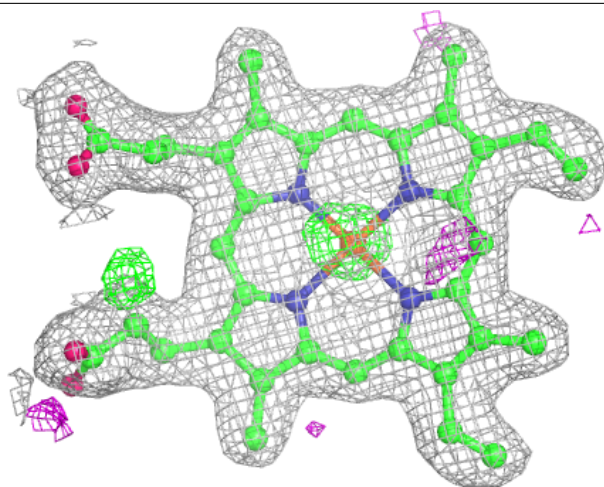
**Electron density around 9P7 A 503:**

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



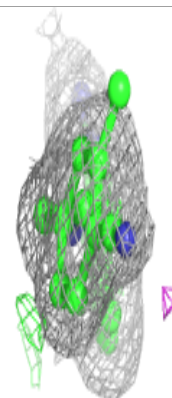
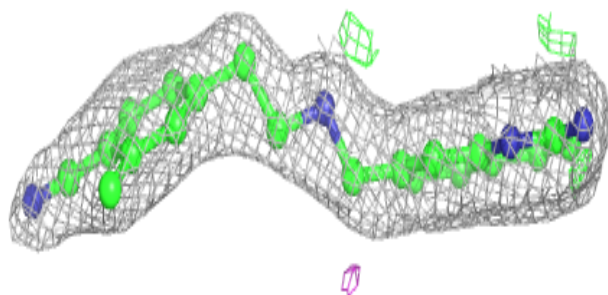
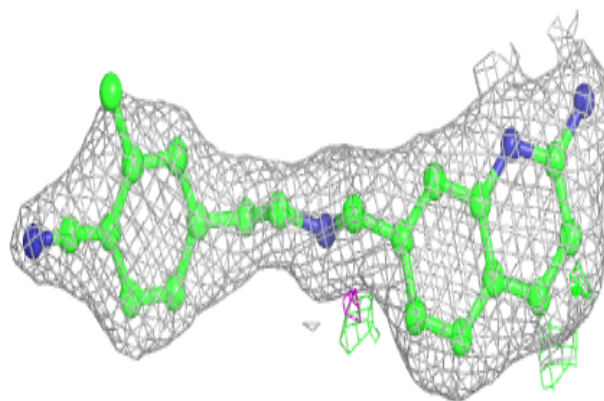
Electron density around HEM B 501:

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

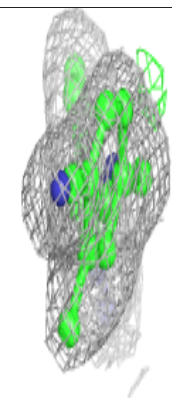
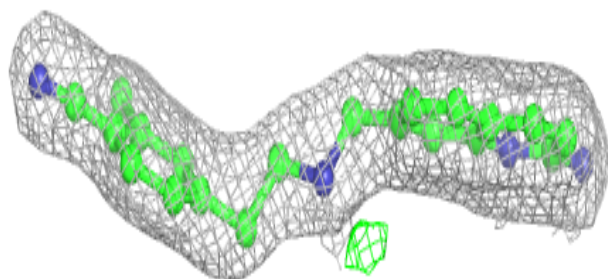
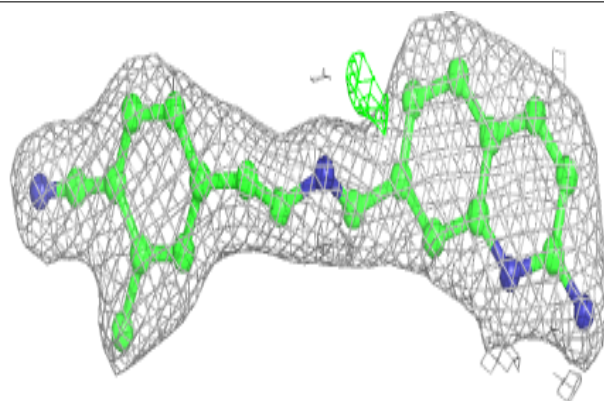


Electron density around 9P7 B 503:

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and green (positive)

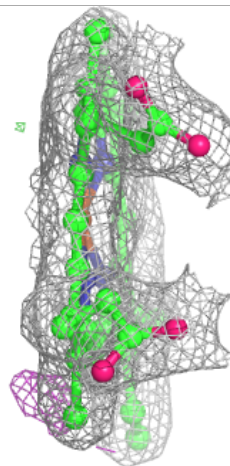
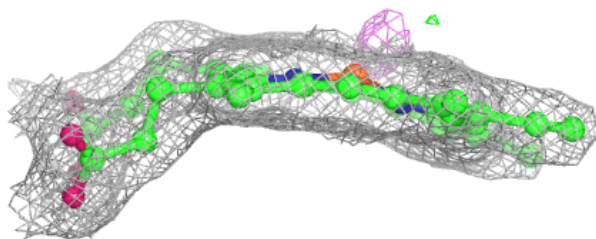
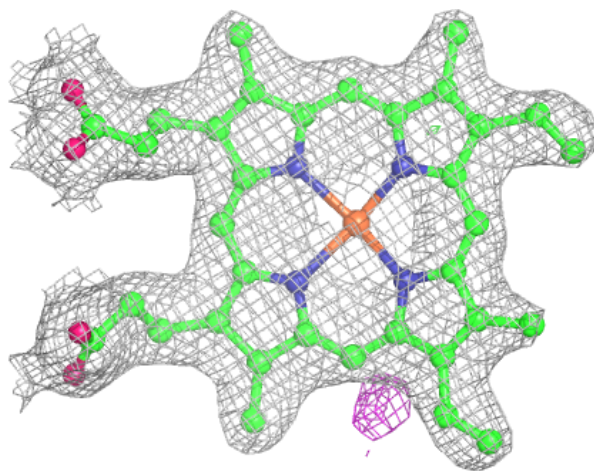
**Electron density around 9P7 D 503:**

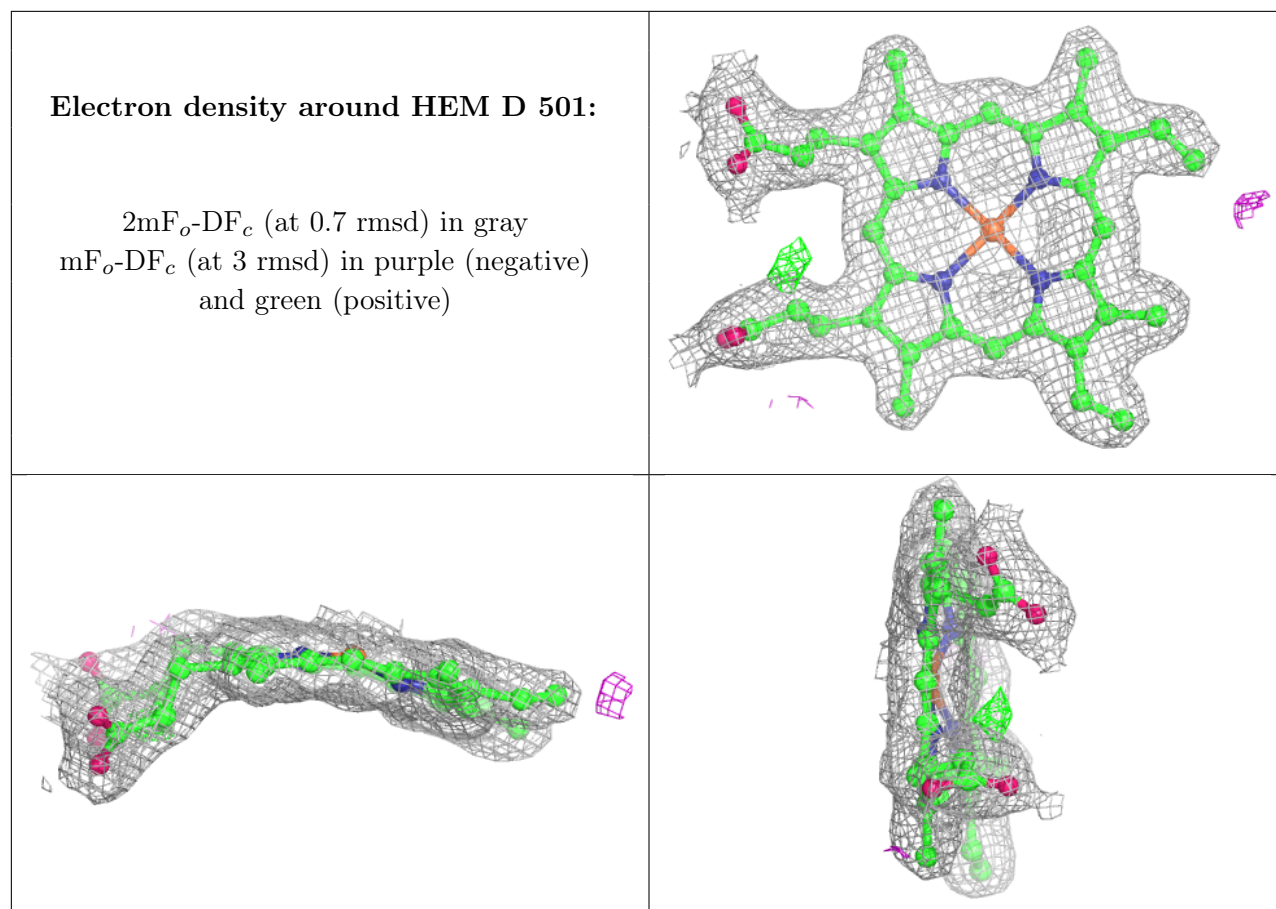
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around HEM A 501:

$2mF_o-DF_c$ (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.