



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 29, 2026 – 09:37 PM UTC

PDB ID : 7ZXD / pdb_00007zxd
Title : Crystal structure of CYP125 from Mycobacterium tuberculosis in complex with an inhibitor
Authors : Snee, M.; Katariya, M.; Levy, C.; Leys, D.
Deposited on : 2022-05-20
Resolution : 2.09 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

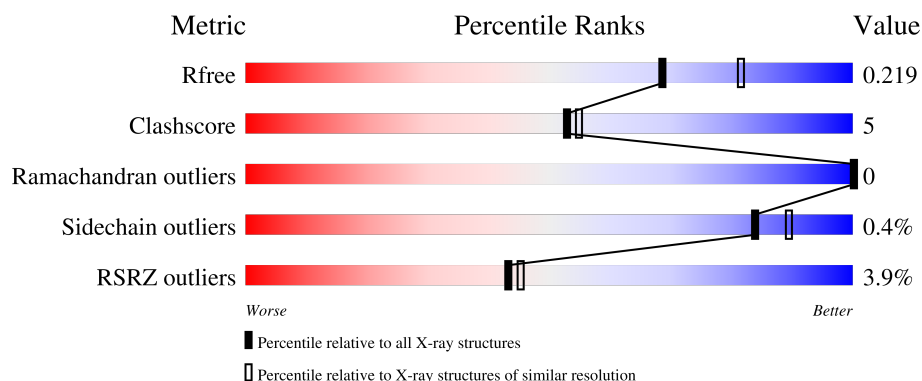
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.09 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}	180053	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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	418	3% 91% 7%
1	B	418	5% 87% 11%
1	C	418	4% 83% 15%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 10380 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.

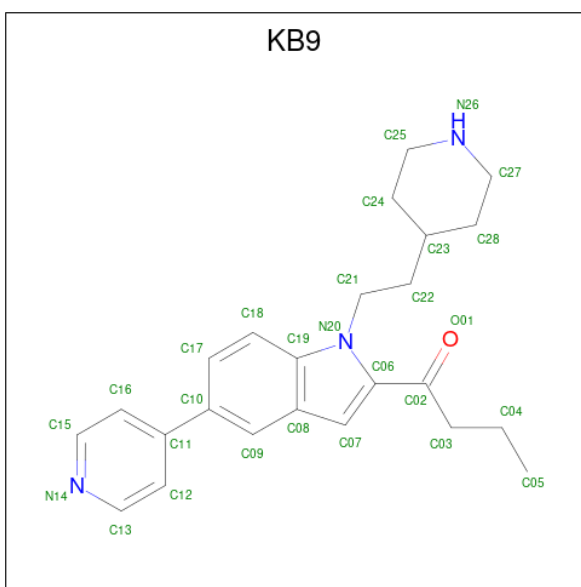
- Molecule 1 is a protein called Steroid C26-monooxygenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	411	Total	C	N	O	S	0	8	0
			3276	2077	578	607	14			
1	B	409	Total	C	N	O	S	0	4	0
			3242	2051	567	610	14			
1	C	409	Total	C	N	O	S	0	5	0
			3232	2046	563	609	14			

There are 6 discrepancies between the modelled and reference sequences:

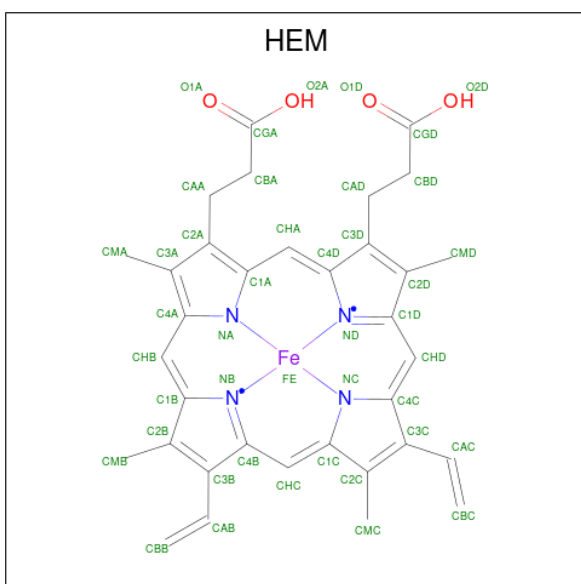
Chain	Residue	Modelled	Actual	Comment	Reference
A	16	ASN	-	expression tag	UNP P9WPP1
A	17	GLY	-	expression tag	UNP P9WPP1
B	16	ASN	-	expression tag	UNP P9WPP1
B	17	GLY	-	expression tag	UNP P9WPP1
C	16	ASN	-	expression tag	UNP P9WPP1
C	17	GLY	-	expression tag	UNP P9WPP1

- Molecule 2 is 1-[1-(2-piperidin-4-ylethyl)-5-pyridin-4-yl-indol-2-yl]butan-1-one (CCD ID: KB9) (formula: C₂₄H₂₉N₃O) (labeled as "Ligand of Interest" by depositor).



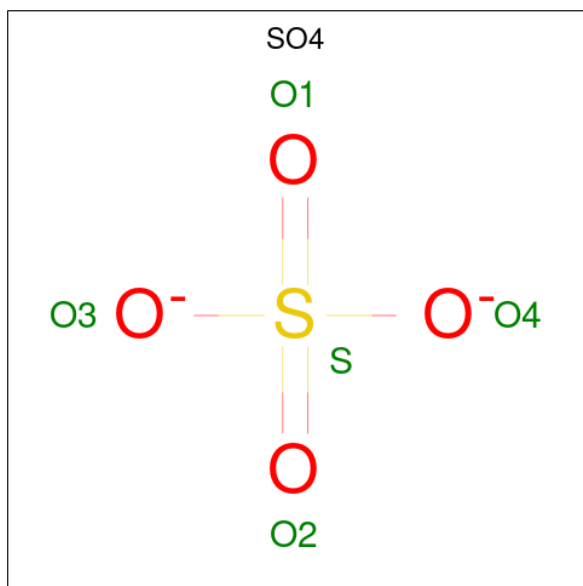
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			28	24	3	1		
2	A	1	Total	C	N	O	0	0
			28	24	3	1		
2	B	1	Total	C	N	O	0	0
			28	24	3	1		
2	C	1	Total	C	N	O	0	0
			28	24	3	1		

- 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	C	Fe	N	O	0	0
			43	34	1	4	4		
3	B	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
3	C	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		

- Molecule 4 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	S	0	0
			5	4	1		
4	A	1	Total	O	S	0	0
			5	4	1		
4	A	1	Total	O	S	0	0
			5	4	1		

- Molecule 5 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

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

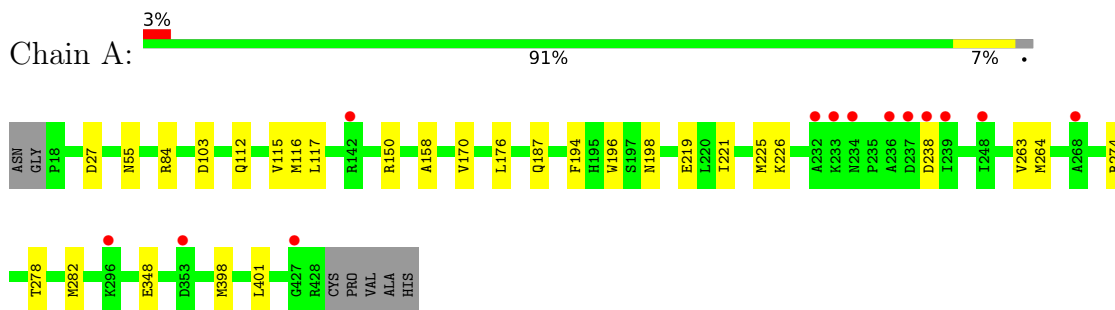
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	228	Total 228	O 228	0	0
6	B	85	Total 85	O 85	0	0
6	C	60	Total 60	O 60	0	0

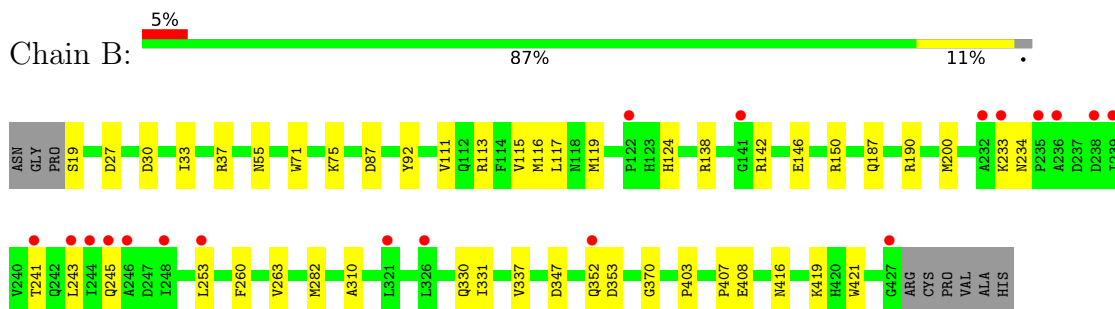
3 Residue-property plots [i](#)

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

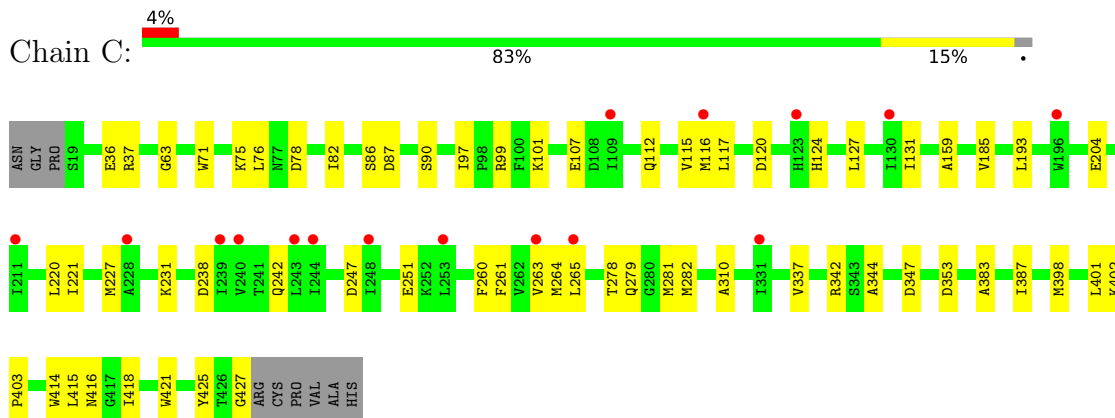
- Molecule 1: Steroid C26-monooxygenase



- Molecule 1: Steroid C26-monooxygenase



- Molecule 1: Steroid C26-monooxygenase



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	137.24Å 69.22Å 144.14Å 90.00° 94.35° 90.00°	Depositor
Resolution (Å)	47.91 – 2.09 47.91 – 2.09	Depositor EDS
% Data completeness (in resolution range)	99.3 (47.91-2.09) 96.9 (47.91-2.09)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.06 (at 2.08Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.196 , 0.219 0.196 , 0.219	Depositor DCC
R_{free} test set	3938 reflections (4.92%)	wwPDB-VP
Wilson B-factor (Å ²)	35.6	Xtriage
Anisotropy	0.225	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 48.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	10380	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.29% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: CL, SO4, KB9, HEM

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	0.09	0/3385	0.29	0/4588
1	B	0.10	0/3329	0.27	0/4520
1	C	0.10	0/3330	0.26	0/4523
All	All	0.10	0/10044	0.27	0/13631

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

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	3276	0	3178	23	0
1	B	3242	0	3097	30	0
1	C	3232	0	3094	42	0
2	A	56	0	0	1	0
2	B	28	0	0	1	0
2	C	28	0	0	2	0
3	A	43	0	30	0	0
3	B	43	0	30	4	0
3	C	43	0	30	2	0
4	A	15	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	A	1	0	0	0	0
6	A	228	0	0	2	0
6	B	85	0	0	2	0
6	C	60	0	0	2	0
All	All	10380	0	9459	98	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 (98) 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:B:37:ARG:HD2	1:C:37:ARG:HD2	1.52	0.90
1:C:282:MET:HE1	1:C:418:ILE:HG13	1.58	0.85
3:C:502:HEM:HBC2	3:C:502:HEM:HHD	1.69	0.75
1:C:281:MET:HE2	1:C:401:LEU:HD22	1.70	0.73
1:A:150[B]:ARG:HH11	1:A:150[B]:ARG:HB2	1.56	0.71
1:A:84:ARG:NH2	4:A:505:SO4:O4	2.26	0.68
1:C:63:GLY:HA3	1:C:99:ARG:HD2	1.75	0.68
1:A:112[B]:GLN:OE1	6:A:601:HOH:O	2.15	0.65
1:C:279:GLN:HA	1:C:282:MET:HE3	1.79	0.65
1:C:278:THR:HG22	1:C:282:MET:HE2	1.77	0.64
1:A:116:MET:HE2	1:A:264:MET:HE1	1.79	0.64
1:A:103:ASP:OD1	6:A:602:HOH:O	2.15	0.63
1:C:99:ARG:NH1	6:C:602:HOH:O	2.32	0.63
1:C:159:ALA:HB2	1:C:398:MET:HG2	1.83	0.61
3:B:502:HEM:HBB2	3:B:502:HEM:HMB1	1.82	0.60
1:C:116:MET:H	1:C:264:MET:HE2	1.69	0.58
1:C:115:VAL:HB	1:C:264:MET:HG3	1.87	0.56
1:C:36:GLU:OE1	6:C:601:HOH:O	2.18	0.56
1:C:107:GLU:N	1:C:107:GLU:OE1	2.39	0.56
1:B:115:VAL:C	1:B:119:MET:HE3	2.32	0.55
1:B:233:LYS:C	1:B:234:ASN:HD22	2.14	0.55
1:C:402:LYS:HD2	1:C:427:GLY:HA2	1.90	0.54
1:C:220:LEU:HB2	1:C:263:VAL:HG12	1.91	0.53
1:A:27:ASP:HA	1:A:55:ASN:HD21	1.74	0.52
1:A:221:ILE:O	1:A:225:MET:HG3	2.10	0.52
1:A:398[B]:MET:HE3	1:A:401:LEU:HD12	1.91	0.52
1:A:117:LEU:HD22	2:A:502:KB9:C16	2.41	0.51
1:C:117:LEU:HD22	2:C:501:KB9:C16	2.41	0.51
3:C:502:HEM:HMB2	3:C:502:HEM:HBB2	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:502:HEM:HMC1	3:B:502:HEM:HBC2	1.92	0.50
1:B:27:ASP:HA	1:B:55:ASN:HD21	1.75	0.49
1:C:185:VAL:HG21	1:C:193:LEU:HD22	1.94	0.48
1:A:150[A]:ARG:HH21	1:A:187:GLN:HE22	1.61	0.48
1:B:71:TRP:CD1	1:B:331:ILE:HD12	2.48	0.48
1:B:116:MET:HG2	1:B:124:HIS:CD2	2.49	0.48
1:C:238:ASP:O	1:C:242:GLN:HG2	2.13	0.48
1:C:247:ASP:OD1	1:C:251:GLU:N	2.44	0.48
1:C:101:LYS:HG2	1:C:204:GLU:HG2	1.95	0.48
1:C:279:GLN:HE22	1:C:416:ASN:ND2	2.12	0.47
1:B:146:GLU:O	1:B:150:ARG:HG3	2.14	0.47
1:B:92:TYR:CZ	1:B:113:ARG:HD3	2.50	0.47
1:B:75:LYS:NZ	1:B:347:ASP:OD1	2.47	0.47
1:B:30:ASP:HB3	1:B:33:ILE:HD12	1.97	0.46
1:B:403:PRO:HB3	1:B:421:TRP:CZ2	2.50	0.46
1:A:226[B]:LYS:HB2	1:A:226[B]:LYS:HE2	1.57	0.46
1:A:348:GLU:H	1:A:348:GLU:CD	2.23	0.46
1:B:243:LEU:HB3	1:B:253:LEU:HD12	1.96	0.46
1:B:408:GLU:HG2	1:B:419:LYS:HE3	1.97	0.46
1:C:97[A]:ILE:HD11	2:C:501:KB9:C18	2.45	0.46
1:A:196:TRP:NE1	1:A:219[A]:GLU:OE1	2.49	0.46
1:C:353:ASP:N	1:C:353:ASP:OD1	2.49	0.46
1:A:170:VAL:HG21	1:A:398[B]:MET:HE1	1.97	0.46
1:B:117:LEU:HD22	2:B:501:KB9:C12	2.46	0.46
1:C:310:ALA:O	1:C:342:ARG:NH2	2.50	0.45
1:A:198:ASN:ND2	1:A:274:ARG:HH22	2.15	0.45
1:C:265:LEU:HD23	1:C:265:LEU:HA	1.78	0.45
1:C:260:PHE:HA	1:C:263:VAL:HG22	1.97	0.45
1:A:158:ALA:HB3	1:A:398[A]:MET:HE1	1.98	0.45
1:A:226[B]:LYS:NZ	1:B:353[B]:ASP:OD2	2.50	0.44
1:B:353[B]:ASP:N	1:B:353[B]:ASP:OD1	2.48	0.44
1:B:200:MET:HE3	1:B:200:MET:HB3	1.81	0.44
1:C:87:ASP:N	1:C:87:ASP:OD1	2.50	0.44
1:C:131:ILE:HD12	1:C:261:PHE:CD1	2.52	0.44
1:B:138:ARG:O	1:B:142:ARG:HB2	2.17	0.44
1:B:241:THR:O	1:B:245:GLN:HG3	2.18	0.44
1:A:150[B]:ARG:HH11	1:A:150[B]:ARG:CB	2.27	0.43
1:B:71:TRP:HB2	1:B:337:VAL:HG12	2.01	0.43
1:B:260:PHE:HA	1:B:263:VAL:HG12	2.00	0.43
1:B:282:MET:HE3	1:B:407:PRO:HG3	2.01	0.43
1:C:78:ASP:O	1:C:82:ILE:HG13	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:403:PRO:HB3	1:C:421:TRP:CZ2	2.54	0.43
1:C:71:TRP:HB2	1:C:337:VAL:HG12	2.00	0.43
1:B:92:TYR:CE2	1:B:113:ARG:HD3	2.54	0.42
1:C:116:MET:HG2	1:C:124:HIS:CD2	2.54	0.42
1:A:176:LEU:HB3	1:A:194:PHE:CE1	2.54	0.42
1:B:419:LYS:NZ	6:B:606:HOH:O	2.48	0.42
3:B:502:HEM:O2A	6:B:601:HOH:O	2.21	0.42
1:A:238:ASP:OD1	1:A:238:ASP:N	2.52	0.42
1:A:278:THR:O	1:A:282:MET:HG3	2.19	0.42
1:C:414:TRP:CZ3	1:C:415:LEU:HD13	2.54	0.42
1:C:90:SER:OG	1:C:120:ASP:OD2	2.28	0.42
1:C:127:LEU:O	1:C:131:ILE:HG12	2.20	0.42
1:B:310:ALA:O	1:B:416:ASN:ND2	2.52	0.42
1:C:221:ILE:HG13	1:C:263:VAL:HG11	2.02	0.42
1:C:76:LEU:HD12	1:C:344:ALA:O	2.20	0.41
1:B:370:GLY:HA3	3:B:502:HEM:HBA1	2.02	0.41
1:C:401:LEU:HD12	1:C:425:TYR:HD1	1.85	0.41
1:C:383:ALA:O	1:C:387:ILE:HG13	2.20	0.41
1:B:87:ASP:OD1	1:B:87:ASP:N	2.52	0.41
1:A:221:ILE:HG13	1:A:263:VAL:HG21	2.02	0.41
1:C:75:LYS:NZ	1:C:347:ASP:OD1	2.53	0.41
1:B:330:GLN:H	1:B:330:GLN:CD	2.26	0.41
1:C:112:GLN:O	1:C:115:VAL:HG22	2.20	0.41
1:B:187:GLN:HA	1:B:190:ARG:HG3	2.03	0.41
1:C:116:MET:HE2	1:C:116:MET:HB3	1.81	0.41
1:B:55:ASN:HD22	1:B:55:ASN:C	2.29	0.40
1:C:227:MET:O	1:C:231:LYS:HG2	2.21	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	417/418 (100%)	407 (98%)	10 (2%)	0	100	100
1	B	411/418 (98%)	401 (98%)	10 (2%)	0	100	100
1	C	412/418 (99%)	404 (98%)	8 (2%)	0	100	100
All	All	1240/1254 (99%)	1212 (98%)	28 (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	341/341 (100%)	341 (100%)	0	100	100
1	B	335/341 (98%)	331 (99%)	4 (1%)	63	72
1	C	335/341 (98%)	334 (100%)	1 (0%)	86	91
All	All	1011/1023 (99%)	1006 (100%)	5 (0%)	84	88

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	19	SER
1	B	111	VAL
1	B	352[A]	GLN
1	B	352[B]	GLN
1	C	86	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	66	HIS
1	A	198	ASN
1	A	234	ASN
1	A	330	GLN
1	B	112	GLN

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Mol	Chain	Res	Type
1	B	148	GLN
1	B	234	ASN
1	B	245	GLN
1	B	375	HIS
1	C	144	HIS
1	C	210	HIS
1	C	242	GLN
1	C	422	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 11 ligands modelled in this entry, 1 is monoatomic - leaving 10 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	SO4	A	504	-	4,4,4	0.24	0	6,6,6	0.07	0
2	KB9	B	501	-	30,31,31	1.24	3 (10%)	40,42,42	1.70	7 (17%)
3	HEM	C	502	1,6	50,50,50	1.43	9 (18%)	67,82,82	1.21	7 (10%)
4	SO4	A	506	-	4,4,4	0.24	0	6,6,6	0.08	0
3	HEM	B	502	1,6	50,50,50	1.39	9 (18%)	67,82,82	1.12	4 (5%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	KB9	A	502	-	30,31,31	1.25	3 (10%)	40,42,42	1.43	4 (10%)
4	SO4	A	505	-	4,4,4	0.24	0	6,6,6	0.07	0
2	KB9	A	501	-	30,31,31	1.24	3 (10%)	40,42,42	1.45	3 (7%)
2	KB9	C	501	-	30,31,31	1.27	3 (10%)	40,42,42	1.43	4 (10%)
3	HEM	A	503	1,6	50,50,50	1.38	8 (16%)	67,82,82	1.09	5 (7%)

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	KB9	B	501	-	-	7/16/24/24	0/4/4/4
3	HEM	C	502	1,6	-	1/14/54/54	-
3	HEM	B	502	1,6	-	0/14/54/54	-
2	KB9	A	502	-	-	5/16/24/24	0/4/4/4
2	KB9	A	501	-	-	1/16/24/24	0/4/4/4
2	KB9	C	501	-	-	4/16/24/24	0/4/4/4
3	HEM	A	503	1,6	-	0/14/54/54	-

All (38) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	501	KB9	C06-N20	-4.31	1.34	1.39
2	B	501	KB9	C06-N20	-4.26	1.34	1.39
2	A	502	KB9	C06-N20	-4.25	1.34	1.39
2	A	501	KB9	C06-N20	-4.21	1.34	1.39
3	C	502	HEM	FE-NB	3.86	2.06	1.94
3	A	503	HEM	FE-NA	3.30	2.06	1.95
3	B	502	HEM	FE-NA	3.16	2.05	1.95
3	C	502	HEM	CAB-C3B	3.10	1.55	1.47
3	B	502	HEM	FE-NC	3.10	2.05	1.95
3	B	502	HEM	CAB-C3B	3.09	1.55	1.47
3	C	502	HEM	FE-NC	3.09	2.05	1.95
3	B	502	HEM	CAC-C3C	3.08	1.55	1.47
3	A	503	HEM	CAC-C3C	3.07	1.55	1.47
3	C	502	HEM	FE-ND	3.04	2.04	1.94
3	B	502	HEM	FE-ND	3.03	2.04	1.94
3	A	503	HEM	CAB-C3B	2.98	1.55	1.47
3	A	503	HEM	FE-ND	2.89	2.03	1.94
3	C	502	HEM	FE-NA	2.84	2.04	1.95

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	502	HEM	FE-NB	2.84	2.03	1.94
3	A	503	HEM	FE-NC	2.83	2.04	1.95
3	A	503	HEM	FE-NB	2.66	2.03	1.94
2	C	501	KB9	C08-C19	-2.65	1.37	1.41
2	A	501	KB9	C08-C19	-2.64	1.37	1.41
3	C	502	HEM	CAC-C3C	2.60	1.54	1.47
2	B	501	KB9	C08-C19	-2.60	1.37	1.41
2	A	502	KB9	C08-C19	-2.53	1.37	1.41
2	A	502	KB9	O01-C02	-2.17	1.18	1.23
3	A	503	HEM	CMC-C2C	2.15	1.55	1.50
3	B	502	HEM	CMC-C2C	2.13	1.55	1.50
2	A	501	KB9	O01-C02	-2.09	1.18	1.23
2	B	501	KB9	O01-C02	-2.07	1.18	1.23
2	C	501	KB9	O01-C02	-2.06	1.18	1.23
3	A	503	HEM	CMB-C2B	2.04	1.55	1.50
3	B	502	HEM	CMB-C2B	2.03	1.54	1.50
3	B	502	HEM	CMA-C3A	2.03	1.54	1.50
3	C	502	HEM	CMA-C3A	2.01	1.54	1.50
3	C	502	HEM	CMC-C2C	2.01	1.54	1.50
3	C	502	HEM	CMD-C2D	2.00	1.54	1.50

All (34) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	KB9	C07-C06-C02	-5.37	124.41	130.49
2	A	502	KB9	C07-C06-C02	-5.21	124.58	130.49
2	B	501	KB9	C07-C06-C02	-4.77	125.09	130.49
2	C	501	KB9	C07-C06-C02	-4.52	125.37	130.49
2	B	501	KB9	C27-C28-C23	4.33	118.84	112.14
2	C	501	KB9	C07-C06-N20	-3.92	107.20	109.42
2	A	501	KB9	C07-C06-N20	-3.59	107.39	109.42
2	B	501	KB9	C07-C06-N20	-3.56	107.41	109.42
2	B	501	KB9	C28-C23-C24	3.42	117.64	109.29
2	A	502	KB9	C07-C06-N20	-2.99	107.73	109.42
3	C	502	HEM	C4D-ND-C1D	2.71	108.42	105.21
2	B	501	KB9	C21-C22-C23	-2.64	109.02	114.85
3	B	502	HEM	C4D-ND-C1D	2.50	108.17	105.21
2	C	501	KB9	C21-C22-C23	-2.48	109.37	114.85
3	A	503	HEM	C4D-ND-C1D	2.46	108.12	105.21
3	B	502	HEM	C1B-NB-C4B	2.45	108.11	105.21
3	B	502	HEM	C3B-C2B-C1B	2.40	108.22	106.41
3	C	502	HEM	C3B-C2B-C1B	2.27	108.11	106.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	501	KB9	C19-N20-C06	2.22	110.07	108.44
3	C	502	HEM	C4C-C3C-C2C	2.22	108.73	106.81
2	B	501	KB9	C28-C27-N26	2.21	114.97	110.81
3	C	502	HEM	C3D-C4D-ND	-2.19	107.77	110.17
3	C	502	HEM	C2A-C1A-NA	-2.17	107.74	110.15
3	B	502	HEM	C3D-C4D-ND	-2.17	107.79	110.17
2	A	502	KB9	C15-N14-C13	2.17	121.83	116.86
2	A	502	KB9	C17-C10-C11	-2.16	117.53	121.25
3	A	503	HEM	C3D-C4D-ND	-2.15	107.81	110.17
3	C	502	HEM	CHC-C1C-NC	2.12	126.76	124.45
3	C	502	HEM	C2D-C1D-ND	-2.11	107.47	109.90
3	A	503	HEM	C2A-C1A-NA	-2.07	107.86	110.15
2	B	501	KB9	C19-N20-C06	2.06	109.95	108.44
3	A	503	HEM	C1B-NB-C4B	2.06	107.65	105.21
3	A	503	HEM	C3B-C2B-C1B	2.04	107.94	106.41
2	A	501	KB9	C15-N14-C13	2.00	121.44	116.86

There are no chirality outliers.

All (18) torsion outliers are listed below:

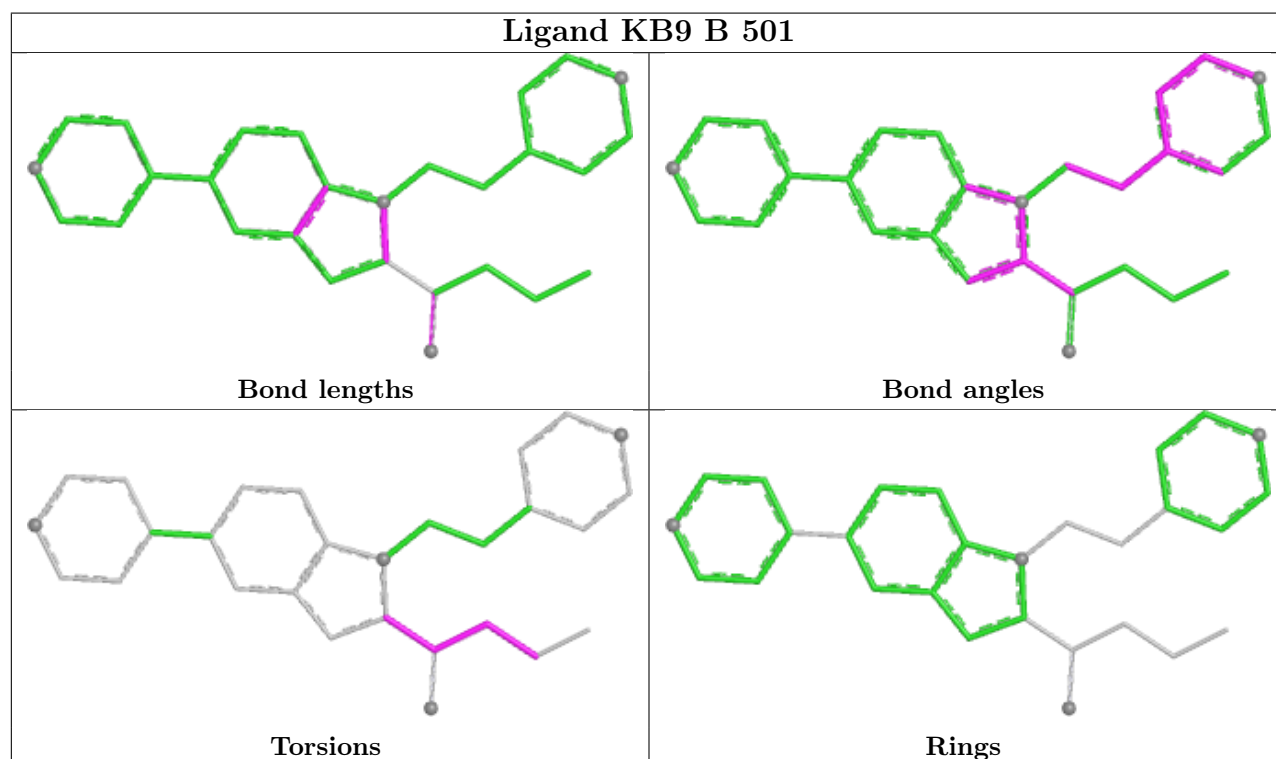
Mol	Chain	Res	Type	Atoms
2	A	502	KB9	O01-C02-C06-N20
2	B	501	KB9	O01-C02-C06-N20
2	C	501	KB9	O01-C02-C06-N20
2	B	501	KB9	C02-C03-C04-C05
2	C	501	KB9	C02-C03-C04-C05
2	A	502	KB9	C21-C22-C23-C28
2	C	501	KB9	C21-C22-C23-C28
3	C	502	HEM	C4C-C3C-CAC-CBC
2	B	501	KB9	C06-C02-C03-C04
2	A	502	KB9	C03-C02-C06-C07
2	A	502	KB9	O01-C02-C06-C07
2	B	501	KB9	C03-C02-C06-C07
2	B	501	KB9	O01-C02-C06-C07
2	C	501	KB9	O01-C02-C06-C07
2	A	501	KB9	O01-C02-C03-C04
2	A	502	KB9	C03-C02-C06-N20
2	B	501	KB9	C03-C02-C06-N20
2	B	501	KB9	O01-C02-C03-C04

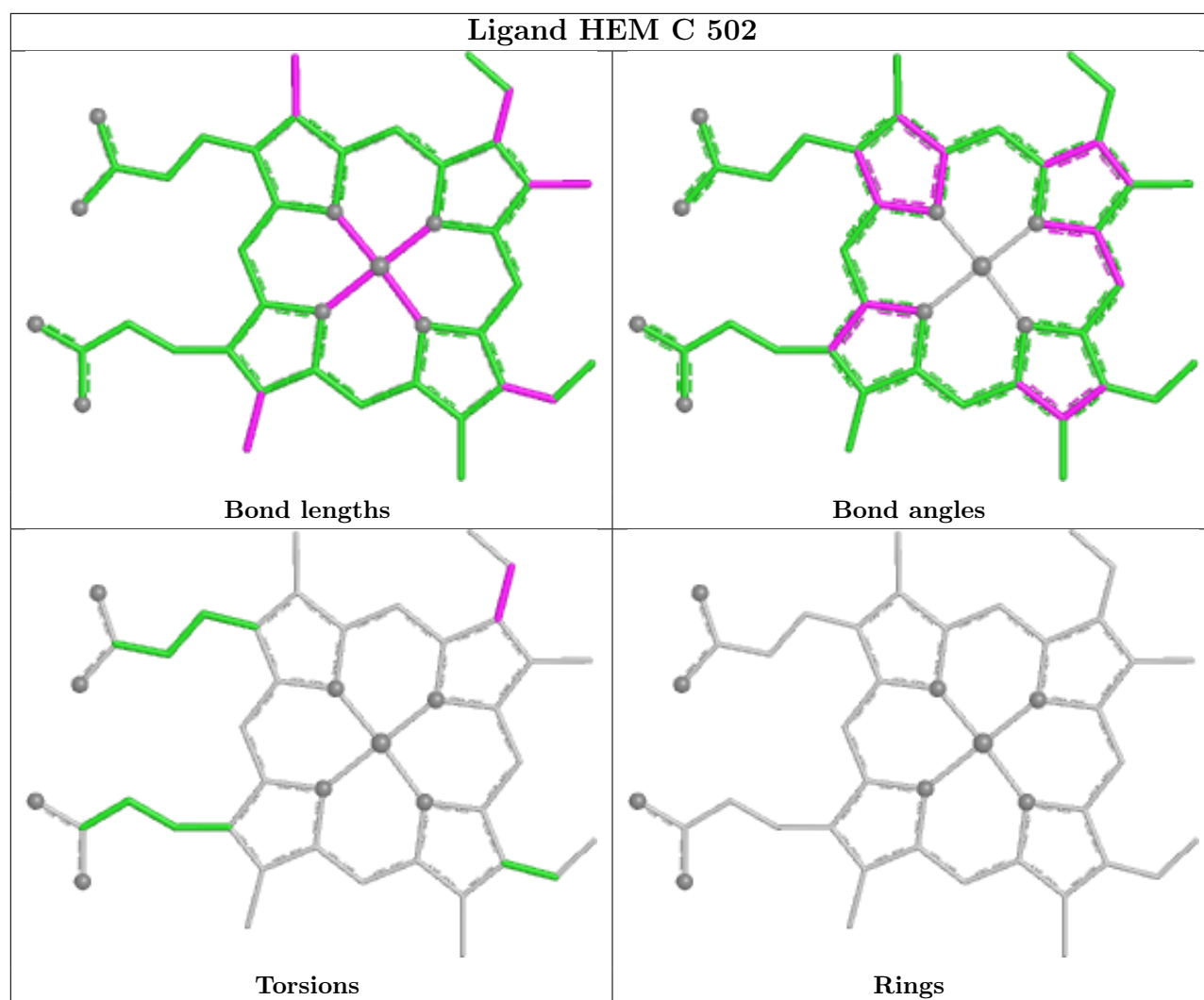
There are no ring outliers.

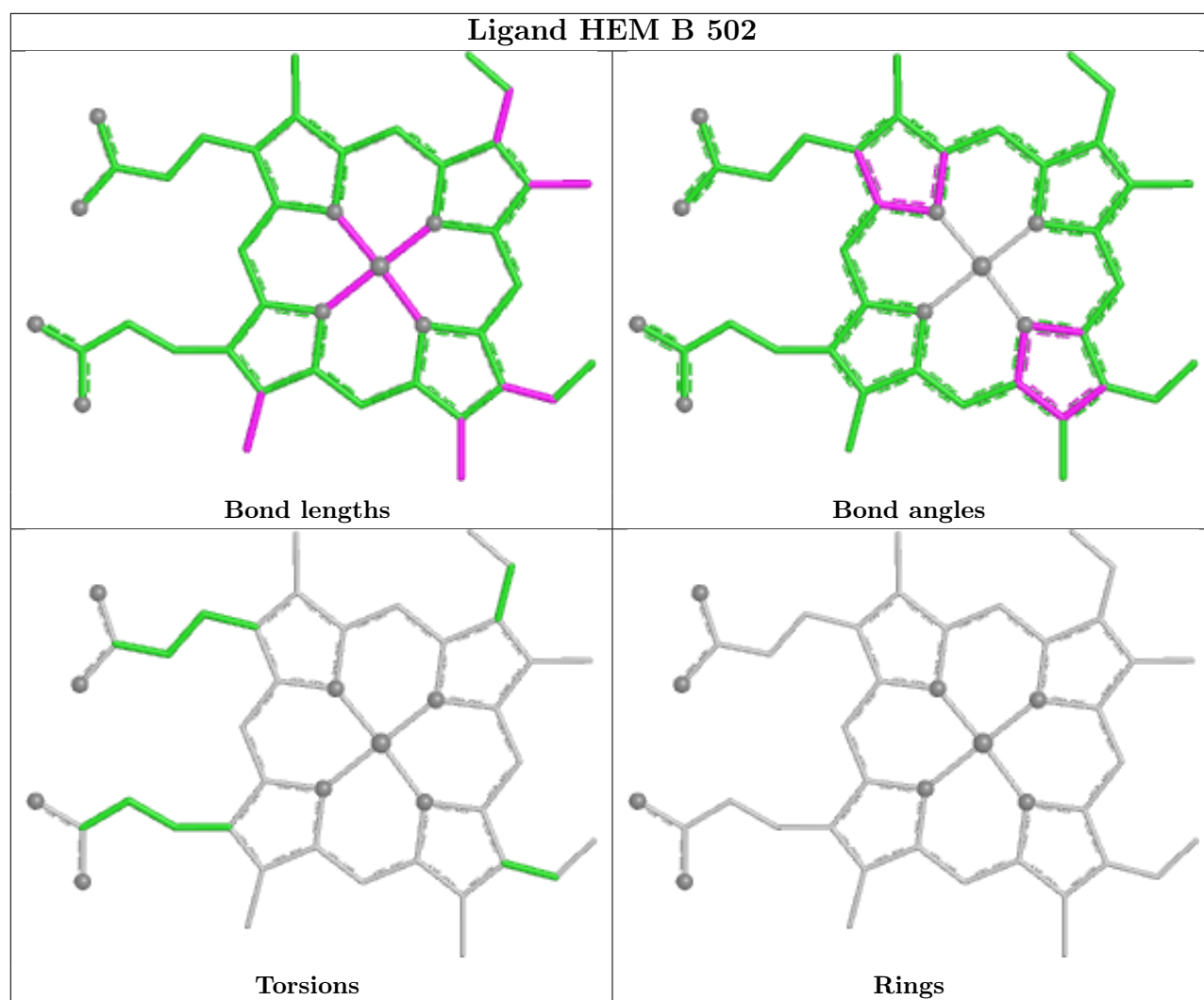
6 monomers are involved in 11 short contacts:

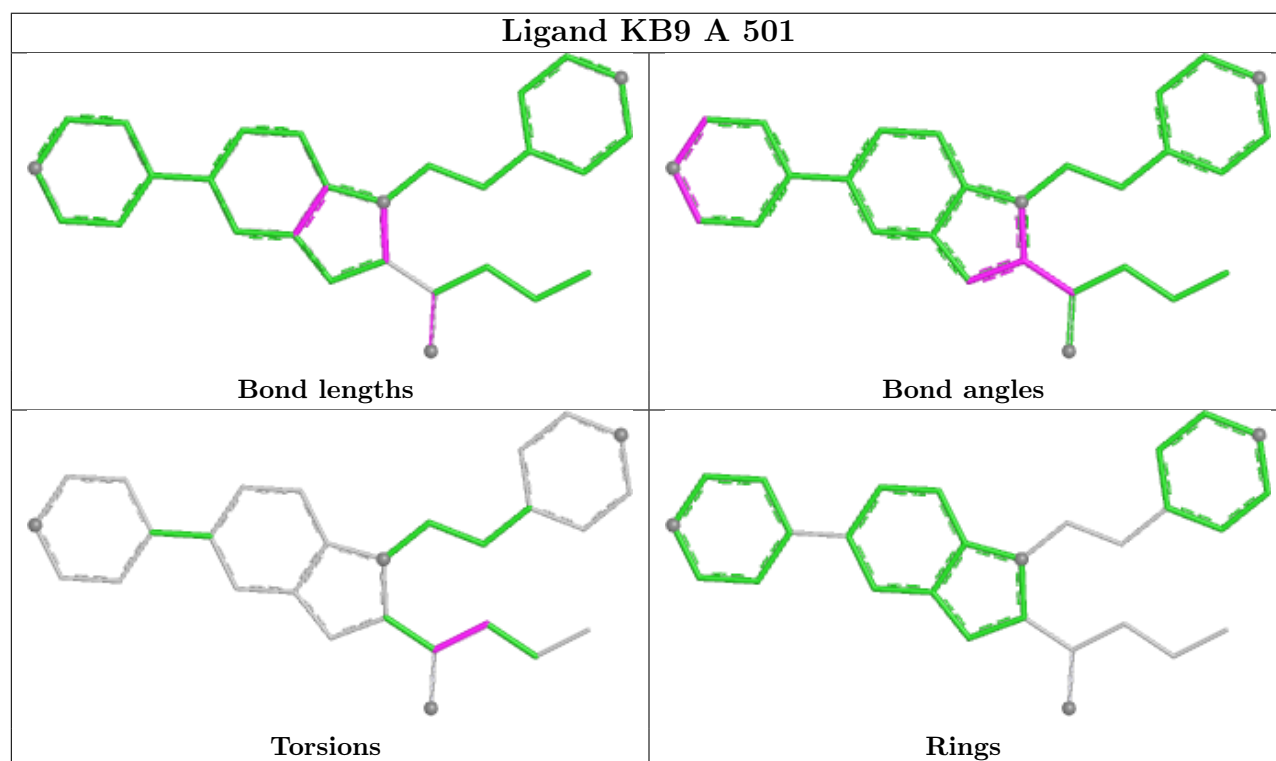
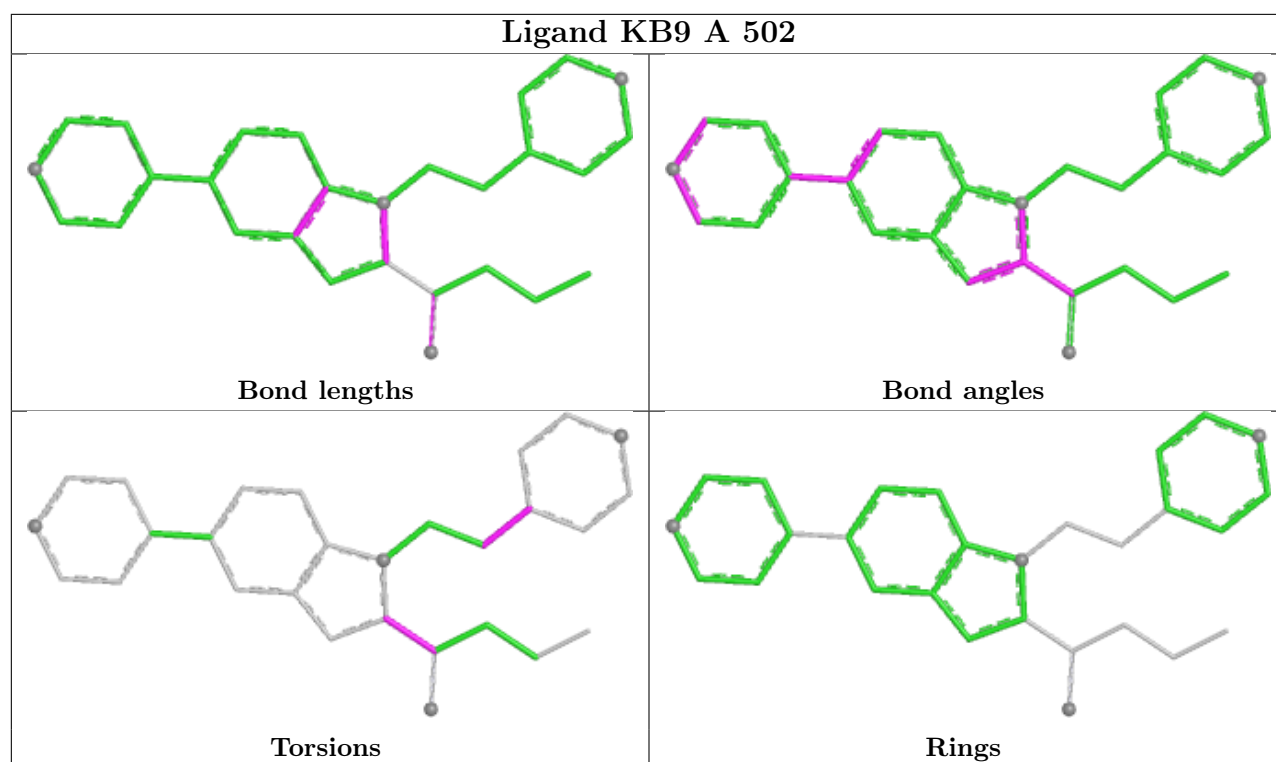
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	501	KB9	1	0
3	C	502	HEM	2	0
3	B	502	HEM	4	0
2	A	502	KB9	1	0
4	A	505	SO4	1	0
2	C	501	KB9	2	0

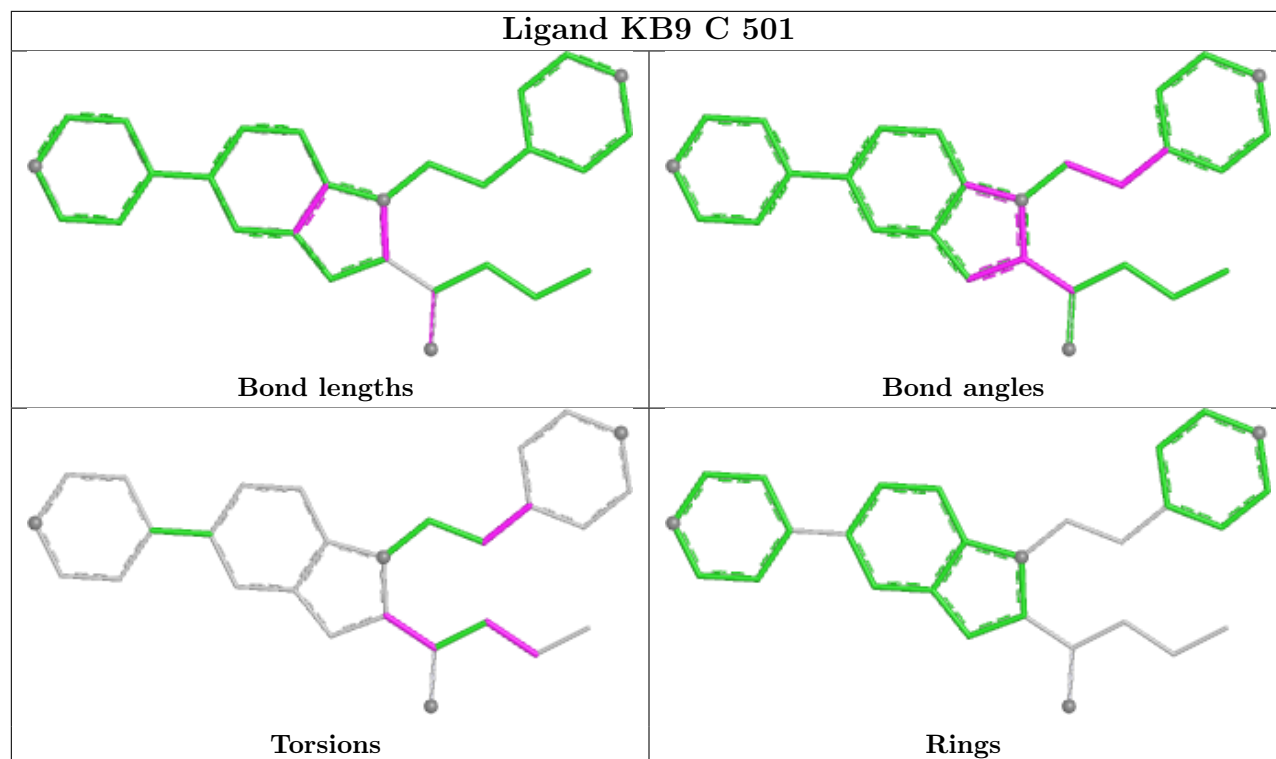
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

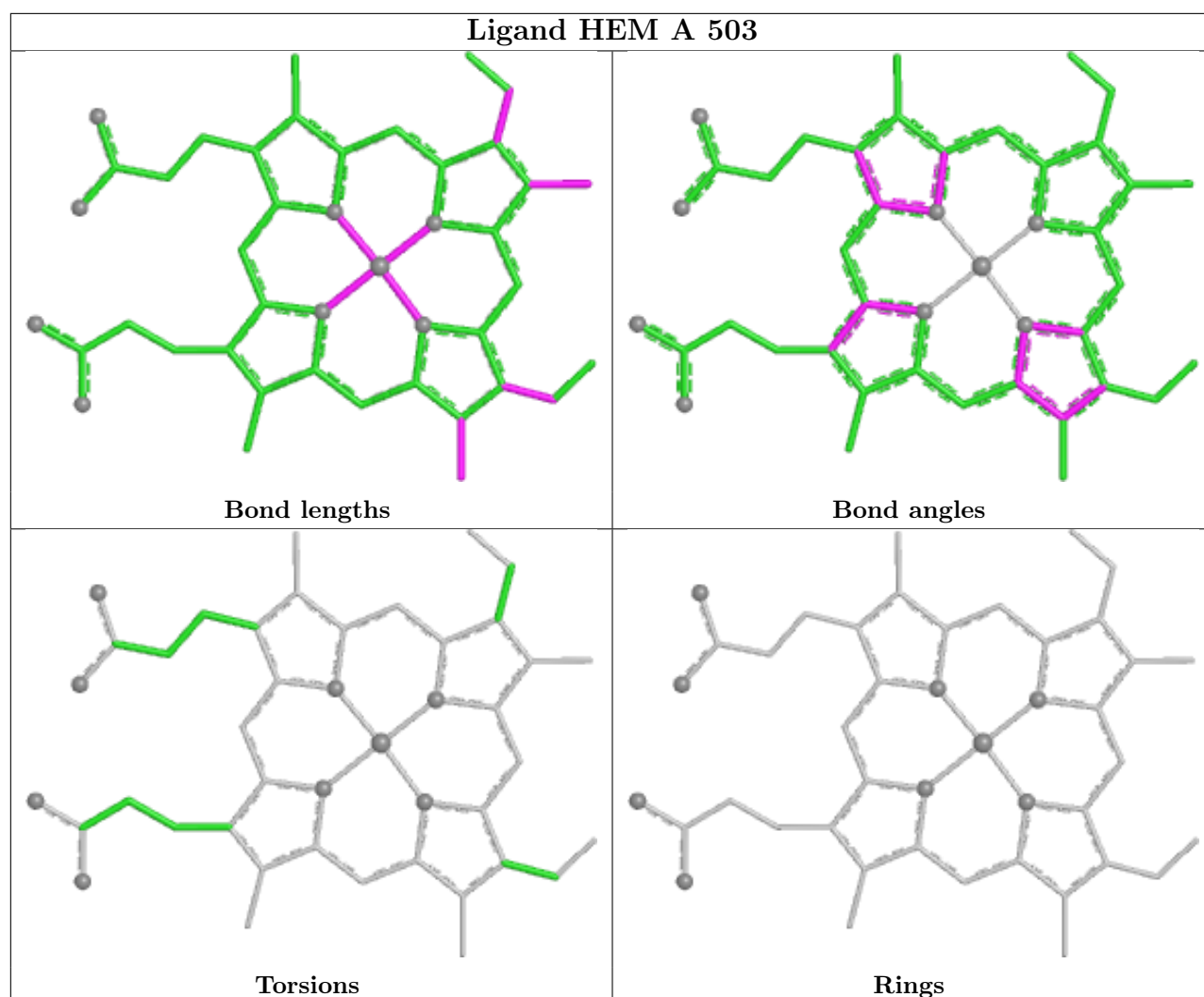












5.7 Other polymers [i](#)

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 ⓘ

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	411/418 (98%)	-0.01	13 (3%)	50	53	16, 33, 56, 101	8 (1%)
1	B	409/418 (97%)	0.55	19 (4%)	37	39	21, 54, 90, 128	4 (0%)
1	C	409/418 (97%)	0.72	16 (3%)	43	45	32, 61, 84, 103	5 (1%)
All	All	1229/1254 (98%)	0.42	48 (3%)	43	45	16, 52, 82, 128	17 (1%)

All (48) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	236	ALA	4.1
1	B	246	ALA	3.7
1	B	248	ILE	3.7
1	A	237	ASP	3.6
1	C	248	ILE	3.4
1	B	235	PRO	3.2
1	A	268	ALA	3.1
1	A	248	ILE	3.1
1	C	130	ILE	3.0
1	B	233	LYS	3.0
1	B	245	GLN	2.9
1	A	236	ALA	2.9
1	A	233	LYS	2.8
1	B	352[A]	GLN	2.8
1	C	211	ILE	2.8
1	B	122	PRO	2.7
1	C	239	ILE	2.7
1	C	196	TRP	2.7
1	A	427	GLY	2.6
1	B	326	LEU	2.6
1	C	265	LEU	2.6
1	B	239	ILE	2.5
1	C	116	MET	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	243	LEU	2.5
1	C	253	LEU	2.5
1	A	238	ASP	2.5
1	A	234	ASN	2.4
1	C	109	ILE	2.4
1	B	241	THR	2.3
1	C	228	ALA	2.3
1	A	353	ASP	2.3
1	A	232	ALA	2.3
1	A	239	ILE	2.2
1	C	263	VAL	2.2
1	B	427	GLY	2.1
1	B	232	ALA	2.1
1	B	253	LEU	2.1
1	B	238	ASP	2.1
1	B	141	GLY	2.1
1	C	244	ILE	2.1
1	B	321	LEU	2.0
1	A	142	ARG	2.0
1	C	123	HIS	2.0
1	B	244	ILE	2.0
1	C	331	ILE	2.0
1	C	240	VAL	2.0
1	C	243	LEU	2.0
1	A	296	LYS	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 ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

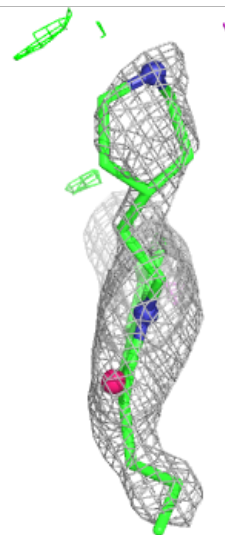
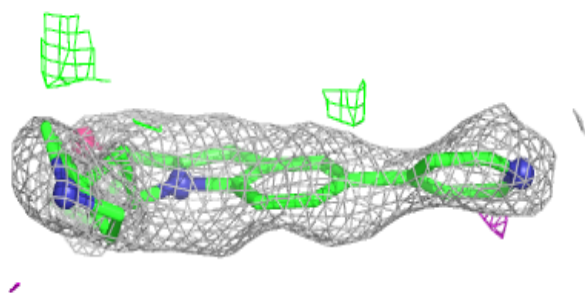
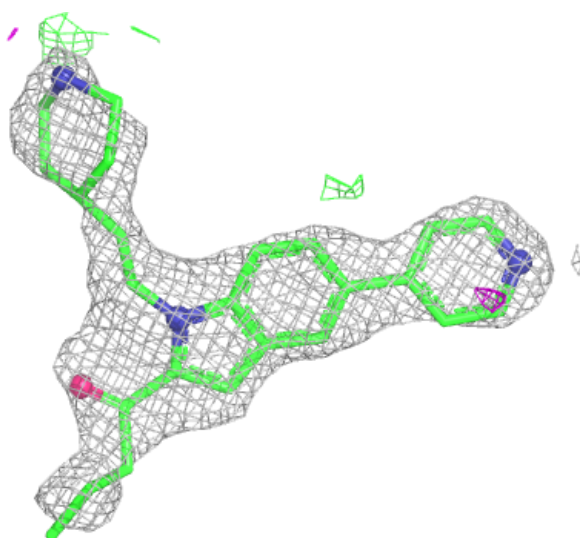
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	SO4	A	505	5/5	0.67	0.11	72,97,114,138	0
4	SO4	A	504	5/5	0.74	0.11	69,74,102,126	0
4	SO4	A	506	5/5	0.83	0.13	44,60,83,86	0
2	KB9	B	501	28/28	0.84	0.16	58,76,91,93	0
2	KB9	C	501	28/28	0.88	0.14	56,67,73,75	0
2	KB9	A	502	28/28	0.90	0.10	26,36,39,41	0
2	KB9	A	501	28/28	0.92	0.10	36,39,46,48	0
3	HEM	C	502	43/43	0.95	0.11	49,54,65,68	0
3	HEM	B	502	43/43	0.97	0.08	40,47,57,60	0
3	HEM	A	503	43/43	0.98	0.06	19,22,28,29	0
5	CL	A	507	1/1	0.98	0.05	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.

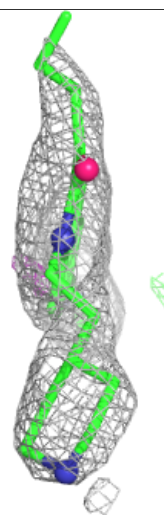
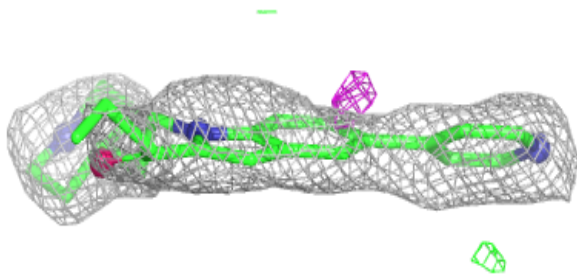
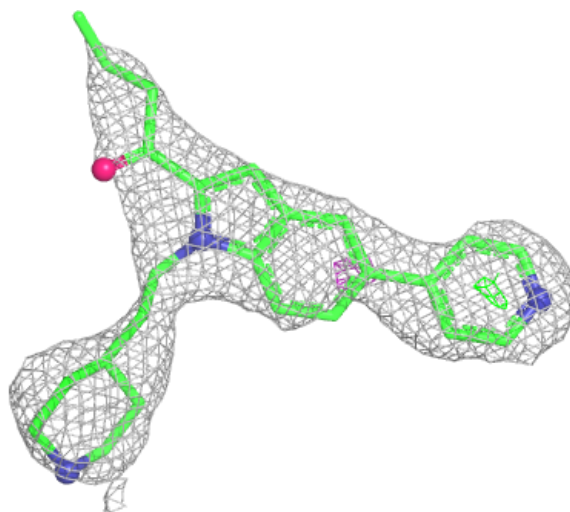
Electron density around KB9 B 501:

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



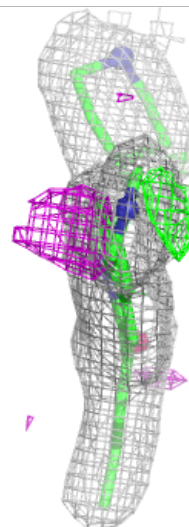
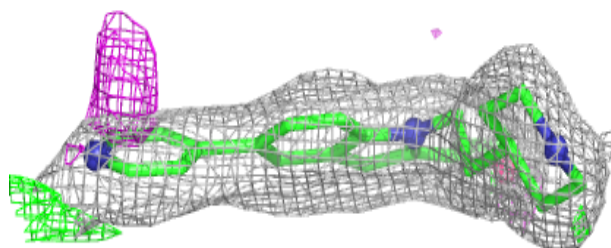
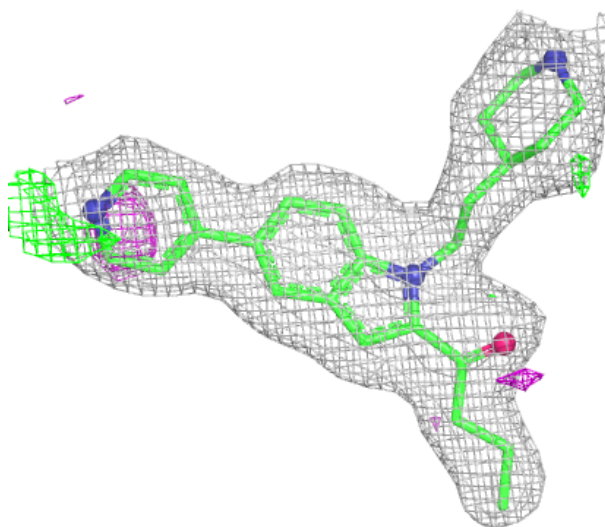
Electron density around KB9 C 501:

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



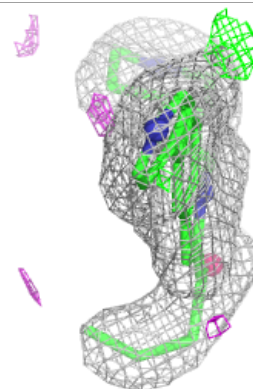
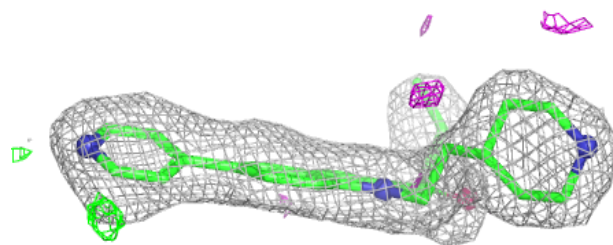
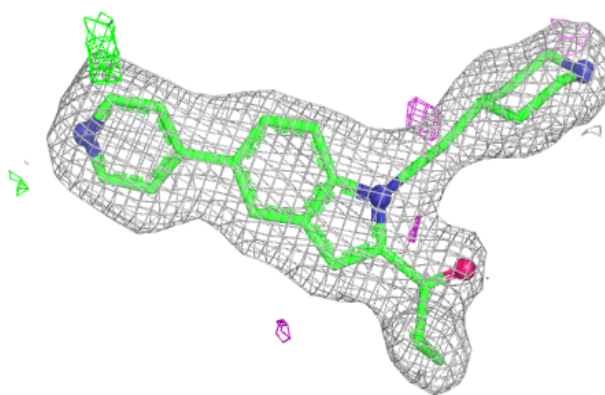
Electron density around KB9 A 502:

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



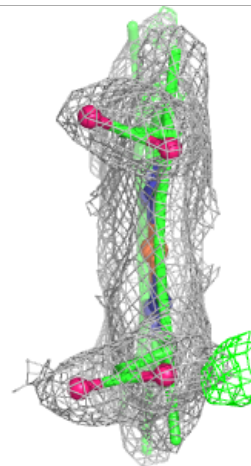
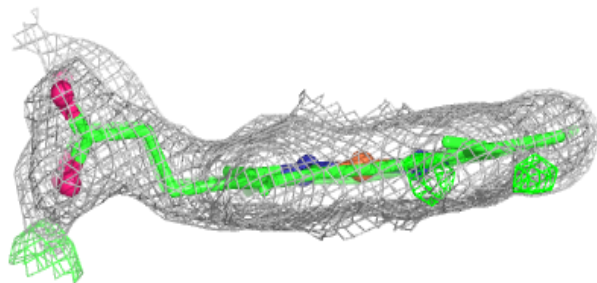
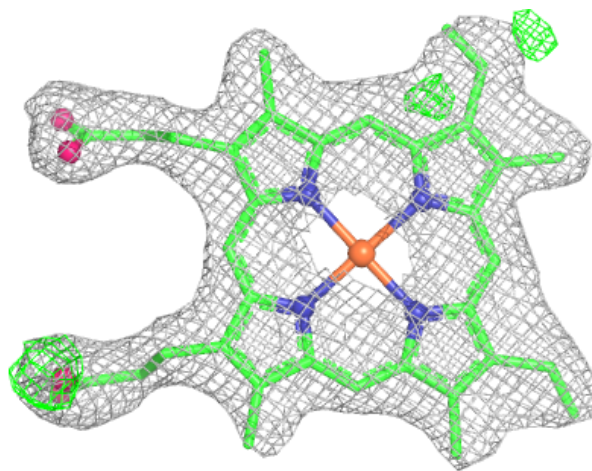
Electron density around KB9 A 501:

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



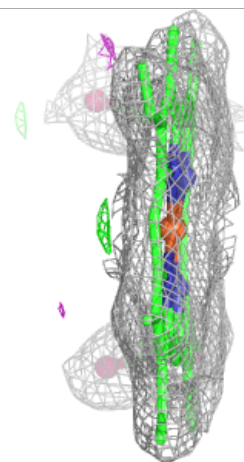
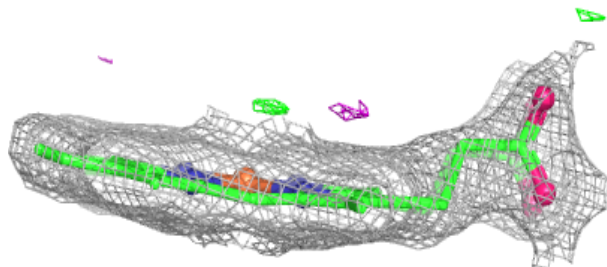
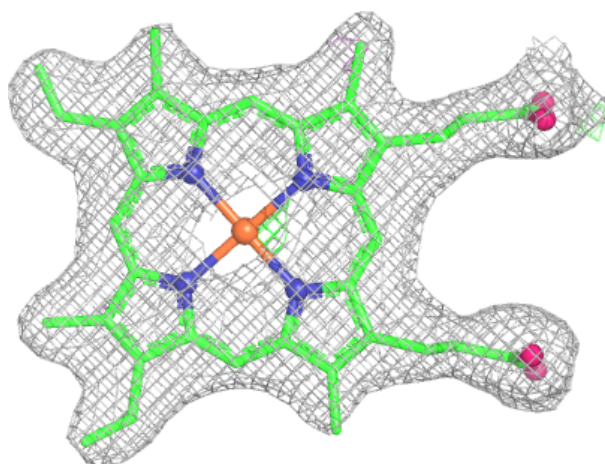
Electron density around HEM C 502:

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



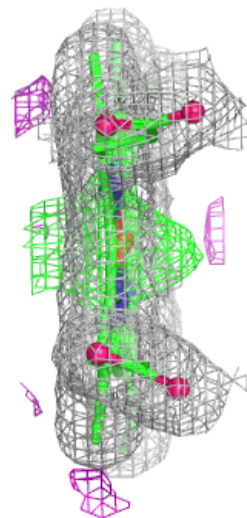
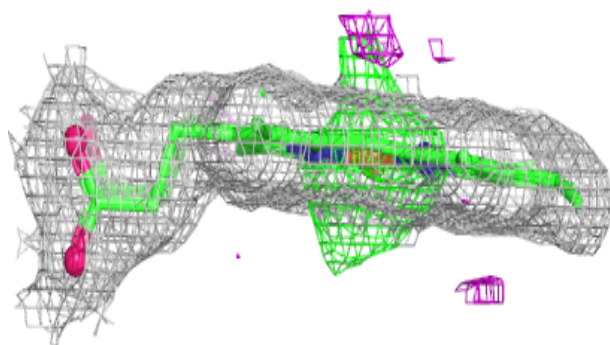
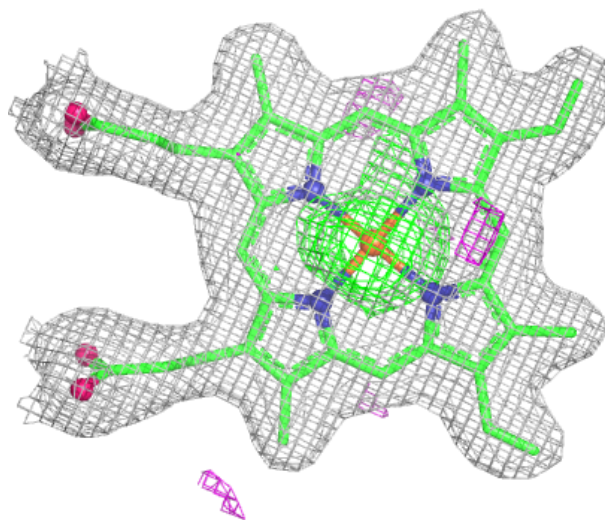
Electron density around HEM B 502:

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



Electron density around HEM A 503:

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



6.5 Other polymers ⓘ

There are no such residues in this entry.