



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 01:57 AM UTC

PDB ID : 6PYY / pdb_00006pyy
Title : Crystal Structure of human Tryptophan 2,3-dioxygenase in complex with PF-06840003 in Active Site and Exo site
Authors : Pham, K.N.; Lewis-Ballester, A.; Yeh, S.R.
Deposited on : 2019-07-31
Resolution : 2.40 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

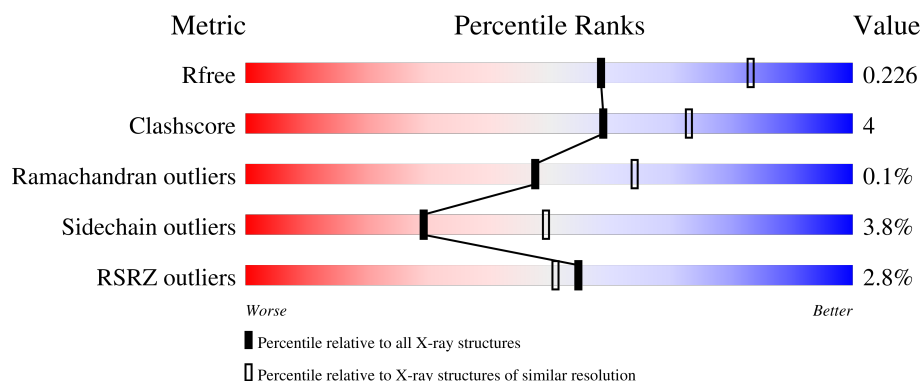
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

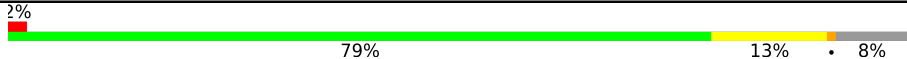
The reported resolution of this entry is 2.40 Å.

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	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

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	380	 2% 79% 13% • 8%
1	B	380	 2% 82% 8% • 10%
1	C	380	 3% 76% 11% • 13%
1	D	380	 2% 82% 8% • 9%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 12156 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 Tryptophan 2,3-dioxygenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	351	Total	C	N	O	S	0	0	0
			2936	1882	516	527	11			
1	B	342	Total	C	N	O	S	0	0	0
			2846	1827	494	514	11			
1	C	332	Total	C	N	O	S	0	0	0
			2749	1770	481	487	11			
1	D	347	Total	C	N	O	S	0	0	0
			2905	1865	513	516	11			

There are 32 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	17	MET	-	initiating methionine	UNP P48775
A	390	GLU	-	expression tag	UNP P48775
A	391	HIS	-	expression tag	UNP P48775
A	392	HIS	-	expression tag	UNP P48775
A	393	HIS	-	expression tag	UNP P48775
A	394	HIS	-	expression tag	UNP P48775
A	395	HIS	-	expression tag	UNP P48775
A	396	HIS	-	expression tag	UNP P48775
B	17	MET	-	initiating methionine	UNP P48775
B	390	GLU	-	expression tag	UNP P48775
B	391	HIS	-	expression tag	UNP P48775
B	392	HIS	-	expression tag	UNP P48775
B	393	HIS	-	expression tag	UNP P48775
B	394	HIS	-	expression tag	UNP P48775
B	395	HIS	-	expression tag	UNP P48775
B	396	HIS	-	expression tag	UNP P48775
C	17	MET	-	initiating methionine	UNP P48775
C	390	GLU	-	expression tag	UNP P48775
C	391	HIS	-	expression tag	UNP P48775
C	392	HIS	-	expression tag	UNP P48775
C	393	HIS	-	expression tag	UNP P48775

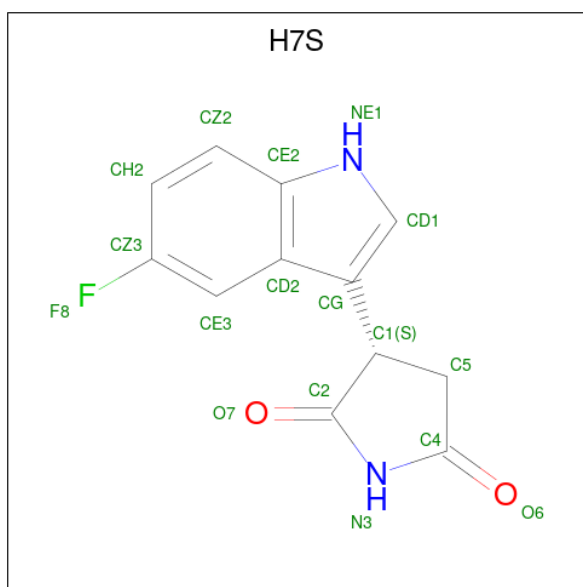
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Chain	Residue	Modelled	Actual	Comment	Reference
C	394	HIS	-	expression tag	UNP P48775
C	395	HIS	-	expression tag	UNP P48775
C	396	HIS	-	expression tag	UNP P48775
D	17	MET	-	initiating methionine	UNP P48775
D	390	GLU	-	expression tag	UNP P48775
D	391	HIS	-	expression tag	UNP P48775
D	392	HIS	-	expression tag	UNP P48775
D	393	HIS	-	expression tag	UNP P48775
D	394	HIS	-	expression tag	UNP P48775
D	395	HIS	-	expression tag	UNP P48775
D	396	HIS	-	expression tag	UNP P48775

- # HEM

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

- 
- WORLD WIDE
PDB
PROTEIN DATA BANK



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	F	N	O	0	0
			17	12	1	2	2		
3	A	1	Total	C	F	N	O	0	0
			17	12	1	2	2		
3	B	1	Total	C	F	N	O	0	0
			17	12	1	2	2		
3	B	1	Total	C	F	N	O	0	0
			17	12	1	2	2		
3	C	1	Total	C	F	N	O	0	0
			17	12	1	2	2		
3	C	1	Total	C	F	N	O	0	0
			17	12	1	2	2		
3	D	1	Total	C	F	N	O	0	0
			17	12	1	2	2		
3	D	1	Total	C	F	N	O	0	0
			17	12	1	2	2		

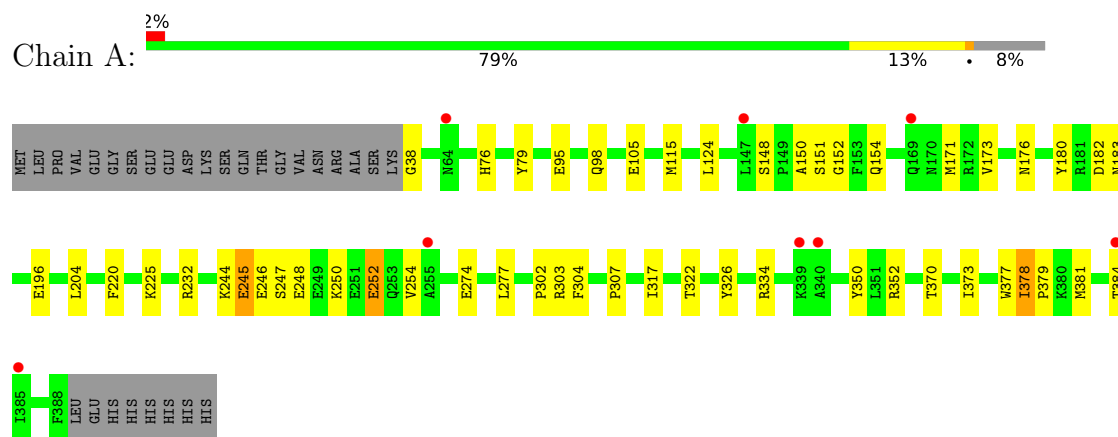
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	114	Total	O	0	0
			114	114		
4	B	111	Total	O	0	0
			111	111		
4	C	80	Total	O	0	0
			80	80		
4	D	107	Total	O	0	0
			107	107		

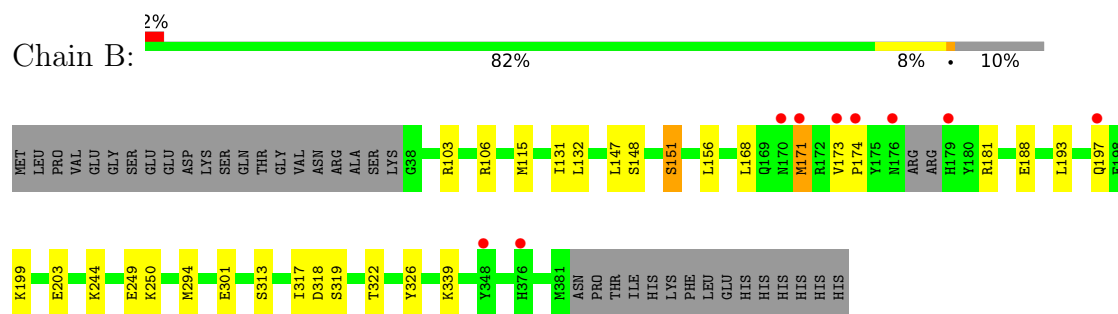
3 Residue-property plots [i](#)

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.

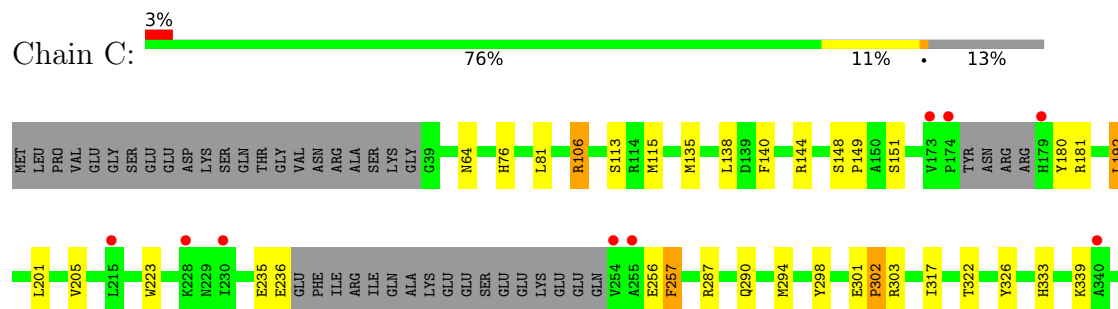
- Molecule 1: Tryptophan 2,3-dioxygenase



- Molecule 1: Tryptophan 2,3-dioxygenase

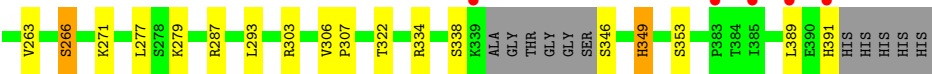
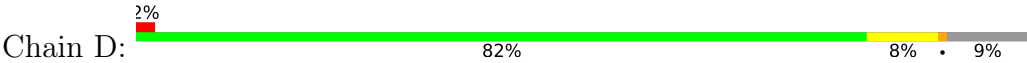


- Molecule 1: Tryptophan 2,3-dioxygenase





● Molecule 1: Tryptophan 2,3-dioxygenase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	144.19Å 154.04Å 87.97Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.90 – 2.40 29.90 – 2.40	Depositor EDS
% Data completeness (in resolution range)	99.8 (29.90-2.40) 99.8 (29.90-2.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.70 (at 2.39Å)	Xtriage
Refinement program	REFMAC 5.8.0238	Depositor
R, R_{free}	0.191 , 0.237 (Not available) , 0.226	Depositor DCC
R_{free} test set	3776 reflections (4.89%)	wwPDB-VP
Wilson B-factor (Å ²)	47.8	Xtriage
Anisotropy	0.067	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 30.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	12156	wwPDB-VP
Average B, all atoms (Å ²)	57.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.93% 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: H7S, 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	1.05	0/3005	1.48	4/4048 (0.1%)
1	B	1.02	0/2910	1.46	3/3918 (0.1%)
1	C	1.05	0/2814	1.47	1/3794 (0.0%)
1	D	1.04	0/2973	1.46	0/4007
All	All	1.04	0/11702	1.47	8/15767 (0.1%)

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	151	SER	CA-C-N	5.56	127.09	120.14
1	B	151	SER	C-N-CA	5.56	127.09	120.14
1	A	38	GLY	CA-C-N	5.54	125.24	119.92
1	A	38	GLY	C-N-CA	5.54	125.24	119.92
1	A	232	ARG	CA-C-N	5.24	125.76	119.94
1	A	232	ARG	C-N-CA	5.24	125.76	119.94
1	C	376	HIS	CA-CB-CG	5.20	119.00	113.80
1	B	318	ASP	CA-CB-CG	5.12	117.72	112.60

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	2936	0	2899	30	0
1	B	2846	0	2808	21	0
1	C	2749	0	2709	31	0
1	D	2905	0	2858	20	0
2	A	43	0	30	6	0
2	B	43	0	30	3	0
2	C	43	0	30	4	0
2	D	43	0	30	5	0
3	A	34	0	0	2	0
3	B	34	0	0	2	0
3	C	34	0	0	1	0
3	D	34	0	0	1	0
4	A	114	0	0	2	0
4	B	111	0	0	2	0
4	C	80	0	0	0	0
4	D	107	0	0	2	0
All	All	12156	0	11394	101	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 4.

All (101) 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:B:401:HEM:HBC2	2:B:401:HEM:HHH	1.74	0.70
2:A:401:HEM:HBB2	2:A:401:HEM:HMB2	1.75	0.69
1:A:150:ALA:HA	1:A:154:GLN:HE22	1.60	0.66
1:C:115:MET:HE3	1:C:317:ILE:CD1	2.29	0.63
1:D:303:ARG:HH22	3:D:403:H7S:C5	2.13	0.61
1:D:346:SER:OG	1:D:349:HIS:HB2	2.00	0.60
1:C:181:ARG:HD3	1:C:192:LEU:HD12	1.84	0.60
2:A:401:HEM:HBC2	2:A:401:HEM:HHH	1.83	0.60
2:C:401:HEM:HMB1	2:C:401:HEM:HBB2	1.84	0.59
1:A:115:MET:HE3	1:A:317:ILE:CD1	2.34	0.57
1:A:245:GLU:O	1:A:247:SER:N	2.36	0.57
2:C:401:HEM:HBC2	2:C:401:HEM:HHH	1.87	0.56
1:B:151:SER:HB2	3:B:402:H7S:O7	2.06	0.55
1:B:115:MET:HE3	1:B:317:ILE:CD1	2.36	0.55
1:B:168:LEU:HD12	1:B:171:MET:HE3	1.87	0.55
1:C:223:TRP:CG	1:C:290:GLN:NE2	2.75	0.55
2:B:401:HEM:HHH	2:B:401:HEM:HBB2	1.89	0.54
1:A:326:TYR:OH	1:C:322:THR:CG2	2.56	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:223:TRP:CD1	1:C:290:GLN:NE2	2.76	0.54
1:C:180:TYR:CE1	1:C:192:LEU:HD11	2.43	0.53
1:C:223:TRP:CD1	1:C:290:GLN:HE21	2.28	0.52
1:D:263:VAL:O	1:D:266:SER:HB2	2.09	0.52
1:D:346:SER:HG	1:D:349:HIS:HB2	1.74	0.52
1:A:176:ASN:ND2	1:A:350:TYR:OH	2.45	0.50
2:D:401:HEM:HBC2	2:D:401:HEM:HHD	1.93	0.50
1:A:151:SER:HB2	3:A:402:H7S:O7	2.10	0.50
1:A:334:ARG:HD2	4:A:559:HOH:O	2.11	0.50
1:C:151:SER:HB2	3:C:402:H7S:O7	2.11	0.50
1:D:263:VAL:O	1:D:266:SER:CB	2.60	0.49
2:C:401:HEM:HBB2	2:C:401:HEM:CMB	2.43	0.49
1:C:140:PHE:CE2	1:C:144:ARG:HD3	2.47	0.49
1:B:326:TYR:OH	1:D:322:THR:HG22	2.12	0.49
1:C:365:LEU:HA	1:C:368:LEU:HD12	1.94	0.49
1:C:76:HIS:NE2	2:C:401:HEM:C2C	2.80	0.49
1:A:150:ALA:HA	1:A:154:GLN:NE2	2.27	0.49
1:A:304:PHE:C	1:A:307:PRO:HD2	2.37	0.49
1:B:301:GLU:OE1	1:C:106:ARG:NH1	2.46	0.48
1:A:76:HIS:NE2	2:A:401:HEM:C2C	2.81	0.48
2:A:401:HEM:HBB2	2:A:401:HEM:CMB	2.43	0.48
1:C:362:PHE:HB3	1:C:365:LEU:HD12	1.96	0.48
1:C:257:PHE:C	1:C:257:PHE:CD1	2.91	0.47
1:D:210:GLU:HG2	1:D:287:ARG:HG2	1.96	0.47
1:B:181:ARG:NH1	1:B:193:LEU:HD11	2.29	0.47
1:C:135:MET:HE3	1:C:140:PHE:HB2	1.97	0.47
1:D:306:VAL:N	1:D:307:PRO:CD	2.78	0.47
2:D:401:HEM:HMB1	2:D:401:HEM:HBB2	1.97	0.47
1:A:105:GLU:OE2	3:A:403:H7S:O6	2.33	0.47
1:C:181:ARG:CD	1:C:192:LEU:HD12	2.44	0.46
1:A:248:GLU:O	1:A:252:GLU:HG2	2.15	0.46
1:C:301:GLU:O	1:C:302:PRO:C	2.59	0.46
1:A:152:GLY:HA3	2:A:401:HEM:C1D	2.50	0.46
1:C:363:VAL:HG22	1:C:367:ASN:ND2	2.30	0.46
1:B:106:ARG:HB3	1:B:106:ARG:CZ	2.46	0.45
1:A:326:TYR:OH	1:C:322:THR:HG22	2.16	0.45
1:B:326:TYR:OH	1:D:322:THR:CG2	2.65	0.45
1:B:339:LYS:NZ	4:B:505:HOH:O	2.50	0.45
2:D:401:HEM:HBB2	2:D:401:HEM:CMB	2.47	0.45
1:A:277:LEU:HD12	1:A:277:LEU:HA	1.89	0.45
1:A:378:ILE:O	1:A:379:PRO:C	2.61	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:223:TRP:CG	1:C:290:GLN:HE21	2.34	0.44
1:B:156:LEU:HD22	1:B:188:GLU:HG2	1.99	0.44
1:A:373:ILE:HD12	1:A:377:TRP:HB2	2.00	0.44
1:D:334:ARG:HD2	4:D:578:HOH:O	2.19	0.43
1:C:180:TYR:HE1	1:C:192:LEU:HD11	1.83	0.43
1:A:303:ARG:HG3	1:D:391:HIS:CB	2.49	0.43
1:C:81:LEU:HD12	1:D:81:LEU:HD12	2.00	0.43
1:C:333:HIS:HB2	1:C:348:TYR:CE1	2.53	0.43
1:C:148:SER:HA	1:C:149:PRO:HA	1.70	0.42
1:B:294:MET:HE2	1:B:294:MET:HB2	1.82	0.42
1:A:182:ASP:OD1	1:A:183:ASN:N	2.52	0.42
1:B:319:SER:O	1:B:322:THR:HG22	2.20	0.42
1:B:244:LYS:HB2	1:B:250:LYS:HD2	2.00	0.42
1:B:132:LEU:CD2	2:B:401:HEM:HBB2	2.49	0.42
1:B:181:ARG:CZ	1:B:193:LEU:HD11	2.50	0.42
1:B:199:LYS:HA	1:B:203:GLU:OE1	2.20	0.42
1:C:303:ARG:HD3	1:C:389:LEU:HA	2.01	0.42
1:A:176:ASN:O	1:A:180:TYR:HB3	2.20	0.42
1:B:173:VAL:O	1:B:174:PRO:C	2.62	0.42
1:D:241:ILE:HD13	1:D:241:ILE:HA	1.89	0.42
1:A:244:LYS:HB2	1:A:250:LYS:HD2	2.01	0.41
1:A:322:THR:HG21	1:C:326:TYR:OH	2.20	0.41
1:D:235:GLU:HA	4:D:592:HOH:O	2.20	0.41
1:A:124:LEU:HD22	1:B:131:ILE:CD1	2.50	0.41
1:B:168:LEU:HB2	1:B:171:MET:HG3	2.02	0.41
1:D:76:HIS:NE2	2:D:401:HEM:C2C	2.89	0.41
1:A:98:GLN:HG3	1:A:204:LEU:HD21	2.03	0.41
1:B:199:LYS:HE2	4:B:605:HOH:O	2.20	0.41
1:C:294:MET:HE3	1:C:298:TYR:CD2	2.56	0.41
1:D:199:LYS:HA	1:D:203:GLU:OE2	2.20	0.41
1:D:293:LEU:HD12	1:D:293:LEU:HA	1.91	0.41
1:A:171:MET:HG3	4:A:579:HOH:O	2.21	0.40
1:C:201:LEU:O	1:C:205:VAL:HG23	2.21	0.40
1:C:287:ARG:O	1:C:290:GLN:HB2	2.21	0.40
1:D:191:LEU:HD12	1:D:191:LEU:HA	1.95	0.40
1:A:79:TYR:CD2	2:A:401:HEM:HAC	2.56	0.40
1:A:302:PRO:HG3	1:D:105:GLU:HG3	2.03	0.40
1:A:326:TYR:OH	1:C:322:THR:HG21	2.21	0.40
1:B:103:ARG:HA	3:B:403:H7S:N3	2.37	0.40
1:A:220:PHE:O	1:A:225:LYS:NZ	2.48	0.40
1:D:152:GLY:HA3	2:D:401:HEM:C1D	2.57	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:370:THR:HA	1:C:339:LYS:HG3	2.04	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	349/380 (92%)	338 (97%)	10 (3%)	1 (0%)	36	50
1	B	338/380 (89%)	328 (97%)	10 (3%)	0	100	100
1	C	326/380 (86%)	302 (93%)	24 (7%)	0	100	100
1	D	336/380 (88%)	327 (97%)	9 (3%)	0	100	100
All	All	1349/1520 (89%)	1295 (96%)	53 (4%)	1 (0%)	48	64

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	246	GLU

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	318/348 (91%)	306 (96%)	12 (4%)	29	49

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	B	307/348 (88%)	301 (98%)	6 (2%)	48 70
1	C	296/348 (85%)	285 (96%)	11 (4%)	30 51
1	D	313/348 (90%)	295 (94%)	18 (6%)	18 32
All	All	1234/1392 (89%)	1187 (96%)	47 (4%)	29 49

All (47) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	95	GLU
1	A	148	SER
1	A	173	VAL
1	A	196	GLU
1	A	245	GLU
1	A	252	GLU
1	A	254	VAL
1	A	274	GLU
1	A	352	ARG
1	A	378	ILE
1	A	381	MET
1	A	384	THR
1	B	147	LEU
1	B	148	SER
1	B	171	MET
1	B	197	GLN
1	B	249	GLU
1	B	313	SER
1	C	64	ASN
1	C	106	ARG
1	C	113	SER
1	C	138	LEU
1	C	192	LEU
1	C	235	GLU
1	C	236	GLU
1	C	256	GLU
1	C	257	PHE
1	C	302	PRO
1	C	375	ARG
1	D	105	GLU
1	D	138	LEU
1	D	148	SER
1	D	149	PRO

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Mol	Chain	Res	Type
1	D	153	PHE
1	D	182	ASP
1	D	194	LYS
1	D	240	ARG
1	D	241	ILE
1	D	249	GLU
1	D	266	SER
1	D	271	LYS
1	D	277	LEU
1	D	279	LYS
1	D	338	SER
1	D	349	HIS
1	D	353	SER
1	D	389	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	116	HIS
1	A	170	ASN
1	A	376	HIS
1	B	163	ASN
1	B	183	ASN
1	B	327	ASN
1	C	44	ASN
1	C	127	GLN
1	C	229	ASN
1	C	290	GLN
1	D	44	ASN
1	D	53	ASN
1	D	55	GLN
1	D	258	GLN
1	D	260	GLN
1	D	290	GLN

5.3.3 RNA ⓘ

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](#)

12 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
2	HEM	D	401	1	50,50,50	1.65	12 (24%)	67,82,82	1.73	17 (25%)
3	H7S	A	403	-	19,19,19	2.10	7 (36%)	23,28,28	2.32	7 (30%)
3	H7S	D	402	-	19,19,19	1.66	3 (15%)	23,28,28	2.64	12 (52%)
2	HEM	C	401	1	50,50,50	1.67	11 (22%)	67,82,82	1.61	13 (19%)
3	H7S	A	402	-	19,19,19	2.01	3 (15%)	23,28,28	2.30	9 (39%)
2	HEM	B	401	1	50,50,50	1.59	9 (18%)	67,82,82	2.08	26 (38%)
3	H7S	C	403	-	19,19,19	2.12	6 (31%)	23,28,28	2.38	8 (34%)
2	HEM	A	401	1	50,50,50	1.62	7 (14%)	67,82,82	1.64	18 (26%)
3	H7S	C	402	-	19,19,19	1.89	3 (15%)	23,28,28	2.35	9 (39%)
3	H7S	D	403	-	19,19,19	1.76	5 (26%)	23,28,28	2.63	10 (43%)
3	H7S	B	403	-	19,19,19	1.86	4 (21%)	23,28,28	2.15	8 (34%)
3	H7S	B	402	-	19,19,19	1.98	4 (21%)	23,28,28	2.39	8 (34%)

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	HEM	D	401	1	-	2/14/54/54	-
3	H7S	A	403	-	-	3/4/16/16	0/3/3/3
3	H7S	D	402	-	-	3/4/16/16	0/3/3/3
2	HEM	C	401	1	-	2/14/54/54	-
3	H7S	A	402	-	-	2/4/16/16	0/3/3/3
2	HEM	B	401	1	-	4/14/54/54	-
3	H7S	C	403	-	-	3/4/16/16	0/3/3/3
2	HEM	A	401	1	-	3/14/54/54	-
3	H7S	C	402	-	-	1/4/16/16	0/3/3/3
3	H7S	D	403	-	-	4/4/16/16	0/3/3/3
3	H7S	B	403	-	-	4/4/16/16	0/3/3/3
3	H7S	B	402	-	-	1/4/16/16	0/3/3/3

All (74) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	403	H7S	CD2-CE2	5.89	1.48	1.41
2	C	401	HEM	FE-NB	5.84	2.12	1.94
3	A	402	H7S	CD2-CE2	5.75	1.48	1.41
3	B	402	H7S	CD2-CE2	5.75	1.48	1.41
2	A	401	HEM	FE-NB	5.42	2.11	1.94
3	C	402	H7S	CD2-CE2	5.23	1.47	1.41
3	D	402	H7S	CD2-CE2	5.19	1.47	1.41
3	A	403	H7S	CD2-CE2	4.96	1.47	1.41
2	B	401	HEM	FE-NB	4.93	2.10	1.94
2	D	401	HEM	FE-NB	4.92	2.10	1.94
3	D	403	H7S	CD2-CE2	4.89	1.47	1.41
3	C	403	H7S	C2-N3	-4.44	1.31	1.37
2	D	401	HEM	FE-NC	4.41	2.09	1.95
3	A	403	H7S	C2-N3	-4.33	1.32	1.37
2	A	401	HEM	FE-NC	4.24	2.09	1.95
2	B	401	HEM	FE-NC	4.19	2.09	1.95
2	C	401	HEM	FE-NC	4.18	2.09	1.95
3	B	403	H7S	C2-N3	-3.94	1.32	1.37
3	A	402	H7S	C4-N3	-3.86	1.31	1.37
2	A	401	HEM	C1B-NB	-3.61	1.34	1.40
3	C	402	H7S	C4-N3	-3.55	1.31	1.37
3	B	403	H7S	CD2-CE2	3.51	1.45	1.41
3	B	403	H7S	CE2-NE1	-3.51	1.32	1.37
3	A	402	H7S	C2-N3	-3.38	1.33	1.37
3	B	402	H7S	C4-N3	-3.33	1.31	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	401	HEM	C4D-ND	-3.32	1.34	1.40
2	B	401	HEM	C4D-ND	-3.29	1.34	1.40
3	A	403	H7S	C4-N3	-3.26	1.32	1.37
2	C	401	HEM	C1C-C2C	-3.20	1.39	1.45
3	B	402	H7S	C2-N3	-3.18	1.33	1.37
2	B	401	HEM	C1B-NB	-3.16	1.34	1.40
3	C	403	H7S	C4-N3	-3.13	1.32	1.37
3	C	402	H7S	C2-N3	-3.02	1.33	1.37
2	A	401	HEM	C1C-C2C	-3.00	1.39	1.45
2	C	401	HEM	FE-NA	2.95	2.04	1.95
2	A	401	HEM	C4B-NB	-2.71	1.33	1.38
2	B	401	HEM	C3C-C4C	-2.65	1.41	1.46
3	D	403	H7S	CD2-CG	2.64	1.49	1.44
2	D	401	HEM	C1B-NB	-2.61	1.35	1.40
2	B	401	HEM	CBD-CAD	2.60	1.60	1.51
3	D	402	H7S	CD1-CG	2.59	1.39	1.36
3	A	403	H7S	CD2-CG	2.57	1.49	1.44
3	D	403	H7S	C4-N3	-2.52	1.33	1.37
3	B	403	H7S	C5-C1	-2.51	1.50	1.54
3	B	402	H7S	C1-CG	-2.50	1.48	1.51
2	C	401	HEM	C1C-NC	-2.49	1.34	1.39
2	C	401	HEM	C3B-C4B	2.47	1.49	1.44
2	D	401	HEM	C3C-C4C	-2.45	1.41	1.46
2	B	401	HEM	CBD-CGD	2.41	1.56	1.50
2	A	401	HEM	O2A-CGA	-2.41	1.22	1.30
3	D	403	H7S	C2-N3	-2.40	1.34	1.37
3	A	403	H7S	C5-C1	-2.37	1.51	1.54
2	D	401	HEM	C3B-C4B	2.33	1.49	1.44
3	A	403	H7S	CE2-NE1	-2.32	1.34	1.37
2	D	401	HEM	C1C-C2C	-2.29	1.40	1.45
2	D	401	HEM	C1A-C2A	-2.26	1.40	1.44
3	C	403	H7S	CD2-CG	2.25	1.48	1.44
3	D	402	H7S	C4-N3	-2.25	1.33	1.37
2	C	401	HEM	C4B-NB	-2.25	1.34	1.38
2	C	401	HEM	C1B-NB	-2.24	1.36	1.40
2	D	401	HEM	C4B-NB	-2.23	1.34	1.38
2	B	401	HEM	C1C-C2C	-2.22	1.40	1.45
2	A	401	HEM	C4D-ND	-2.20	1.36	1.40
2	D	401	HEM	C1D-ND	-2.19	1.34	1.38
3	A	403	H7S	CE3-CD2	-2.17	1.36	1.39
2	B	401	HEM	C3B-C4B	2.16	1.49	1.44
3	C	403	H7S	CE2-NE1	-2.15	1.34	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	401	HEM	C4D-ND	-2.13	1.36	1.40
2	C	401	HEM	C3C-C4C	-2.11	1.42	1.46
2	D	401	HEM	O2D-CGD	-2.06	1.24	1.30
3	C	403	H7S	C1-CG	-2.06	1.49	1.51
2	C	401	HEM	C1D-ND	-2.05	1.34	1.38
2	D	401	HEM	FE-ND	-2.04	1.88	1.94
3	D	403	H7S	CE2-NE1	-2.01	1.34	1.37

All (145) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	403	H7S	C1-CG-CD1	-5.91	118.28	126.75
3	C	403	H7S	C1-CG-CD1	-5.87	118.33	126.75
3	D	403	H7S	C1-CG-CD1	-5.67	118.62	126.75
3	B	402	H7S	CE2-CD2-CG	-5.51	101.85	106.98
3	C	403	H7S	CE2-CD2-CG	-5.41	101.94	106.98
3	B	403	H7S	C1-CG-CD1	-5.27	119.19	126.75
3	D	402	H7S	CZ2-CE2-CD2	-5.25	117.11	122.19
3	C	402	H7S	CE2-CD2-CG	-5.19	102.15	106.98
3	A	402	H7S	CZ2-CE2-CD2	-5.06	117.29	122.19
3	A	402	H7S	CE2-CD2-CG	-4.99	102.33	106.98
2	C	401	HEM	CBD-CAD-C3D	-4.99	98.74	112.53
3	D	403	H7S	CE2-NE1-CD1	4.96	113.49	109.08
3	B	402	H7S	CD2-CG-CD1	4.71	109.85	106.25
3	D	402	H7S	CE2-CD2-CG	-4.71	102.60	106.98
3	D	403	H7S	CE2-CD2-CG	-4.68	102.63	106.98
2	D	401	HEM	CHD-C1D-ND	4.66	129.43	124.42
3	A	402	H7S	CD2-CG-CD1	4.52	109.71	106.25
2	B	401	HEM	C3B-C4B-NB	-4.46	106.27	109.47
2	D	401	HEM	CHC-C4B-NB	4.45	129.21	124.42
3	C	402	H7S	CZ2-CE2-CD2	-4.39	117.94	122.19
2	B	401	HEM	CHB-C1B-NB	4.38	129.78	124.37
3	A	403	H7S	CD2-CG-C1	4.35	134.44	126.56
3	D	402	H7S	CD2-CG-CD1	4.34	109.57	106.25
2	B	401	HEM	C1B-NB-C4B	4.29	110.28	105.21
2	B	401	HEM	CBD-CAD-C3D	-4.12	101.13	112.53
3	B	403	H7S	CD2-CG-C1	4.11	134.01	126.56
3	C	402	H7S	CD2-CG-CD1	4.09	109.38	106.25
3	A	403	H7S	CE2-CD2-CG	-4.08	103.18	106.98
3	B	402	H7S	CZ2-CH2-CZ3	4.08	122.57	118.38
3	D	402	H7S	O6-C4-C5	-3.98	121.41	126.48
3	D	403	H7S	CE3-CD2-CG	3.85	138.80	133.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	401	HEM	CHA-C4D-ND	3.81	129.08	124.37
2	A	401	HEM	CHA-C4D-ND	3.79	129.06	124.37
3	D	402	H7S	O7-C2-C1	-3.77	122.61	126.71
3	C	403	H7S	CE3-CD2-CG	3.73	138.64	133.39
3	A	403	H7S	CE3-CD2-CG	3.73	138.63	133.39
3	D	402	H7S	CE3-CD2-CE2	3.69	122.98	119.39
2	C	401	HEM	CHD-C1D-ND	3.68	128.39	124.42
3	B	403	H7S	CE2-NE1-CD1	3.68	112.35	109.08
2	B	401	HEM	O2D-CGD-CBD	3.61	125.40	114.00
2	C	401	HEM	CHC-C4B-NB	3.60	128.30	124.42
3	B	402	H7S	CZ2-CE2-CD2	-3.59	118.71	122.19
2	D	401	HEM	CBD-CAD-C3D	-3.59	102.60	112.53
2	A	401	HEM	CHC-C4B-NB	3.52	128.21	124.42
3	C	403	H7S	CD2-CG-C1	3.49	132.89	126.56
2	B	401	HEM	CHA-C4D-ND	3.49	128.68	124.37
2	D	401	HEM	CHB-C1B-NB	3.44	128.62	124.37
2	A	401	HEM	CBD-CAD-C3D	-3.43	103.04	112.53
2	D	401	HEM	C1B-NB-C4B	3.42	109.26	105.21
2	B	401	HEM	CHD-C1D-C2D	-3.34	119.75	125.03
2	B	401	HEM	CHD-C1D-ND	3.33	128.01	124.42
3	D	403	H7S	CD2-CG-C1	3.33	132.59	126.56
3	D	403	H7S	O7-C2-C1	-3.33	123.09	126.71
3	C	402	H7S	CE2-NE1-CD1	3.27	111.99	109.08
3	B	403	H7S	CE3-CD2-CG	3.25	137.95	133.39
3	A	403	H7S	C4-N3-C2	-3.24	110.71	113.96
2	B	401	HEM	CHD-C4C-NC	3.23	127.97	124.45
3	D	402	H7S	O6-C4-N3	3.23	129.37	125.06
2	B	401	HEM	C3B-C2B-C1B	3.21	108.83	106.41
3	C	403	H7S	CD2-CG-CD1	3.20	108.70	106.25
2	A	401	HEM	O2A-CGA-O1A	-3.18	115.15	123.33
3	B	403	H7S	CE2-CD2-CG	-3.18	104.02	106.98
3	D	403	H7S	CD2-CG-CD1	3.17	108.67	106.25
2	B	401	HEM	CAB-C3B-C2B	-3.12	118.30	128.43
2	D	401	HEM	CHA-C4D-ND	3.07	128.16	124.37
3	D	403	H7S	CG-CD1-NE1	-3.05	107.83	110.31
3	D	403	H7S	O7-C2-N3	3.00	128.46	124.93
3	A	402	H7S	CE3-CD2-CE2	2.99	122.30	119.39
3	C	402	H7S	CZ2-CH2-CZ3	2.99	121.44	118.38
3	A	403	H7S	CE2-NE1-CD1	2.97	111.72	109.08
2	B	401	HEM	CHC-C4B-NB	2.91	127.56	124.42
3	D	402	H7S	CG-CD1-NE1	-2.91	107.95	110.31
3	A	403	H7S	C5-C4-N3	2.88	111.44	108.02

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	401	HEM	CHD-C1D-ND	2.84	127.48	124.42
2	C	401	HEM	C1B-NB-C4B	2.84	108.56	105.21
3	C	402	H7S	CG-CD1-NE1	-2.82	108.02	110.31
2	D	401	HEM	CAC-C3C-C4C	2.82	131.56	124.82
2	A	401	HEM	C1B-NB-C4B	2.82	108.54	105.21
3	B	402	H7S	C1-CG-CD1	-2.81	122.72	126.75
2	D	401	HEM	CHD-C1D-C2D	-2.78	120.64	125.03
2	B	401	HEM	O2D-CGD-O1D	-2.74	116.29	123.33
3	B	402	H7S	CG-CD1-NE1	-2.73	108.10	110.31
3	C	403	H7S	CE2-NE1-CD1	2.72	111.50	109.08
3	B	402	H7S	CE3-CD2-CE2	2.71	122.03	119.39
3	C	403	H7S	CZ2-CH2-CZ3	2.67	121.12	118.38
3	C	402	H7S	CE3-CD2-CE2	2.66	121.98	119.39
2	A	401	HEM	C4C-NC-C1C	2.66	110.16	105.82
3	A	402	H7S	CE2-NE1-CD1	2.64	111.43	109.08
2	A	401	HEM	O2A-CGA-CBA	2.64	122.34	114.00
3	A	402	H7S	CG-CD1-NE1	-2.63	108.17	110.31
2	B	401	HEM	CHA-C1A-NA	2.62	128.62	123.86
2	A	401	HEM	CHD-C4C-NC	2.62	127.30	124.45
2	B	401	HEM	C1A-CHA-C4D	-2.61	120.12	126.25
2	B	401	HEM	C3C-C2C-C1C	2.61	109.51	107.05
3	B	402	H7S	CE2-NE1-CD1	2.60	111.39	109.08
3	B	403	H7S	O6-C4-C5	-2.59	123.18	126.48
3	A	402	H7S	C5-C4-N3	2.57	111.08	108.02
2	B	401	HEM	C4C-NC-C1C	2.56	110.00	105.82
3	B	403	H7S	C5-C4-N3	2.56	111.06	108.02
2	B	401	HEM	CAB-C3B-C4B	2.53	135.57	124.39
2	A	401	HEM	C1A-CHA-C4D	-2.52	120.33	126.25
2	A	401	HEM	CHB-C1B-NB	2.51	127.47	124.37
2	A	401	HEM	CAC-C3C-C4C	2.50	130.80	124.82
2	D	401	HEM	C3B-C4B-NB	-2.49	107.68	109.47
2	C	401	HEM	O2A-CGA-CBA	2.49	121.86	114.00
2	B	401	HEM	CMB-C2B-C3B	-2.49	122.41	128.43
2	B	401	HEM	O2A-CGA-CBA	2.47	121.81	114.00
2	C	401	HEM	CHB-C1B-NB	2.46	127.41	124.37
2	A	401	HEM	CHA-C4D-C3D	-2.45	120.70	125.23
2	C	401	HEM	O2A-CGA-O1A	-2.45	117.02	123.33
2	B	401	HEM	CMC-C2C-C3C	-2.44	122.52	128.43
3	D	402	H7S	CE2-NE1-CD1	2.41	111.22	109.08
2	D	401	HEM	CHA-C4D-C3D	-2.38	120.84	125.23
2	B	401	HEM	CMB-C2B-C1B	2.38	128.75	125.03
2	C	401	HEM	C1A-CHA-C4D	-2.37	120.68	126.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	401	HEM	CHA-C4D-C3D	-2.32	120.95	125.23
2	D	401	HEM	CAC-C3C-C2C	-2.31	120.92	128.43
2	C	401	HEM	CHD-C1D-C2D	-2.30	121.39	125.03
2	B	401	HEM	CAC-C3C-C4C	2.27	130.24	124.82
3	C	403	H7S	CZ2-CE2-CD2	-2.26	120.00	122.19
3	A	402	H7S	CZ2-CE2-NE1	2.26	134.35	130.74
2	A	401	HEM	C4B-C3B-C2B	-2.24	105.22	107.28
3	C	402	H7S	C5-C4-N3	2.24	110.68	108.02
3	D	403	H7S	CZ2-CE2-CD2	-2.23	120.03	122.19
2	C	401	HEM	C4C-NC-C1C	2.23	109.45	105.82
2	D	401	HEM	CBA-CAA-C2A	-2.22	106.39	112.53
2	C	401	HEM	CAD-C3D-C4D	2.22	128.56	124.70
2	A	401	HEM	CHD-C1D-C2D	-2.21	121.54	125.03
2	A	401	HEM	CAC-C3C-C2C	-2.21	121.26	128.43
2	A	401	HEM	CHC-C1C-NC	2.20	126.85	124.45
3	D	402	H7S	CZ2-CE2-NE1	2.18	134.23	130.74
3	B	403	H7S	C4-N3-C2	-2.17	111.78	113.96
2	D	401	HEM	CAD-C3D-C4D	2.16	128.46	124.70
2	B	401	HEM	C1D-C2D-C3D	-2.15	104.72	106.98
2	D	401	HEM	CHB-C1B-C2B	-2.13	120.88	126.95
2	D	401	HEM	O2D-CGD-CBD	2.13	120.72	114.00
2	B	401	HEM	O2A-CGA-O1A	-2.12	117.87	123.33
2	D	401	HEM	O2A-CGA-O1A	-2.11	117.91	123.33
3	D	402	H7S	O7-C2-N3	2.08	127.38	124.93
2	C	401	HEM	CHA-C4D-C3D	-2.06	121.43	125.23
3	C	402	H7S	O6-C4-N3	-2.04	122.34	125.06
2	D	401	HEM	C1A-CHA-C4D	-2.04	121.45	126.25
3	A	402	H7S	CZ2-CH2-CZ3	2.04	120.47	118.38
2	A	401	HEM	CMD-C2D-C1D	2.03	128.21	125.03
3	D	402	H7S	CZ2-CH2-CZ3	2.02	120.46	118.38

There are no chirality outliers.

All (32) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	403	H7S	C5-C1-CG-CD1
3	A	403	H7S	C5-C1-CG-CD2
3	B	403	H7S	C5-C1-CG-CD1
3	B	403	H7S	C5-C1-CG-CD2
3	C	403	H7S	C5-C1-CG-CD1
3	C	403	H7S	C5-C1-CG-CD2
3	D	403	H7S	C5-C1-CG-CD1

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Mol	Chain	Res	Type	Atoms
3	D	403	H7S	C5-C1-CG-CD2
2	B	401	HEM	C4B-C3B-CAB-CBB
2	B	401	HEM	C4C-C3C-CAC-CBC
3	B	403	H7S	C2-C1-CG-CD2
3	C	403	H7S	C2-C1-CG-CD2
3	C	402	H7S	C5-C1-CG-CD2
2	B	401	HEM	CAA-CBA-CGA-O1A
2	B	401	HEM	CAA-CBA-CGA-O2A
2	D	401	HEM	CAA-CBA-CGA-O1A
2	A	401	HEM	CAA-CBA-CGA-O1A
2	C	401	HEM	CAA-CBA-CGA-O1A
2	C	401	HEM	CAA-CBA-CGA-O2A
2	A	401	HEM	CAA-CBA-CGA-O2A
2	D	401	HEM	CAA-CBA-CGA-O2A
2	A	401	HEM	C4C-C3C-CAC-CBC
3	A	403	H7S	C2-C1-CG-CD2
3	B	402	H7S	C2-C1-CG-CD2
3	D	403	H7S	C2-C1-CG-CD2
3	A	402	H7S	C5-C1-CG-CD1
3	D	402	H7S	C5-C1-CG-CD1
3	B	403	H7S	C2-C1-CG-CD1
3	D	402	H7S	C2-C1-CG-CD1
3	D	403	H7S	C2-C1-CG-CD1
3	A	402	H7S	C5-C1-CG-CD2
3	D	402	H7S	C5-C1-CG-CD2

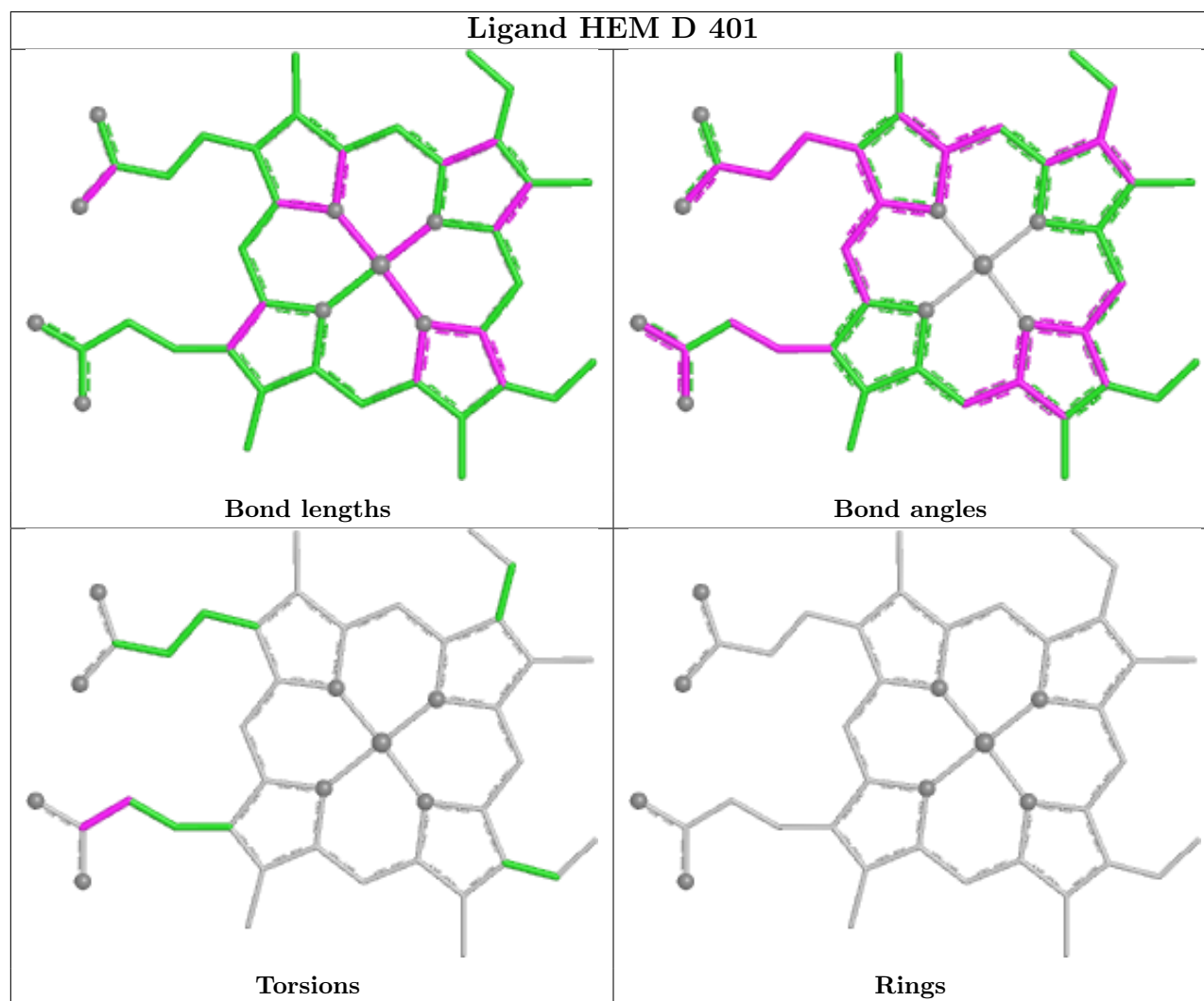
There are no ring outliers.

10 monomers are involved in 24 short contacts:

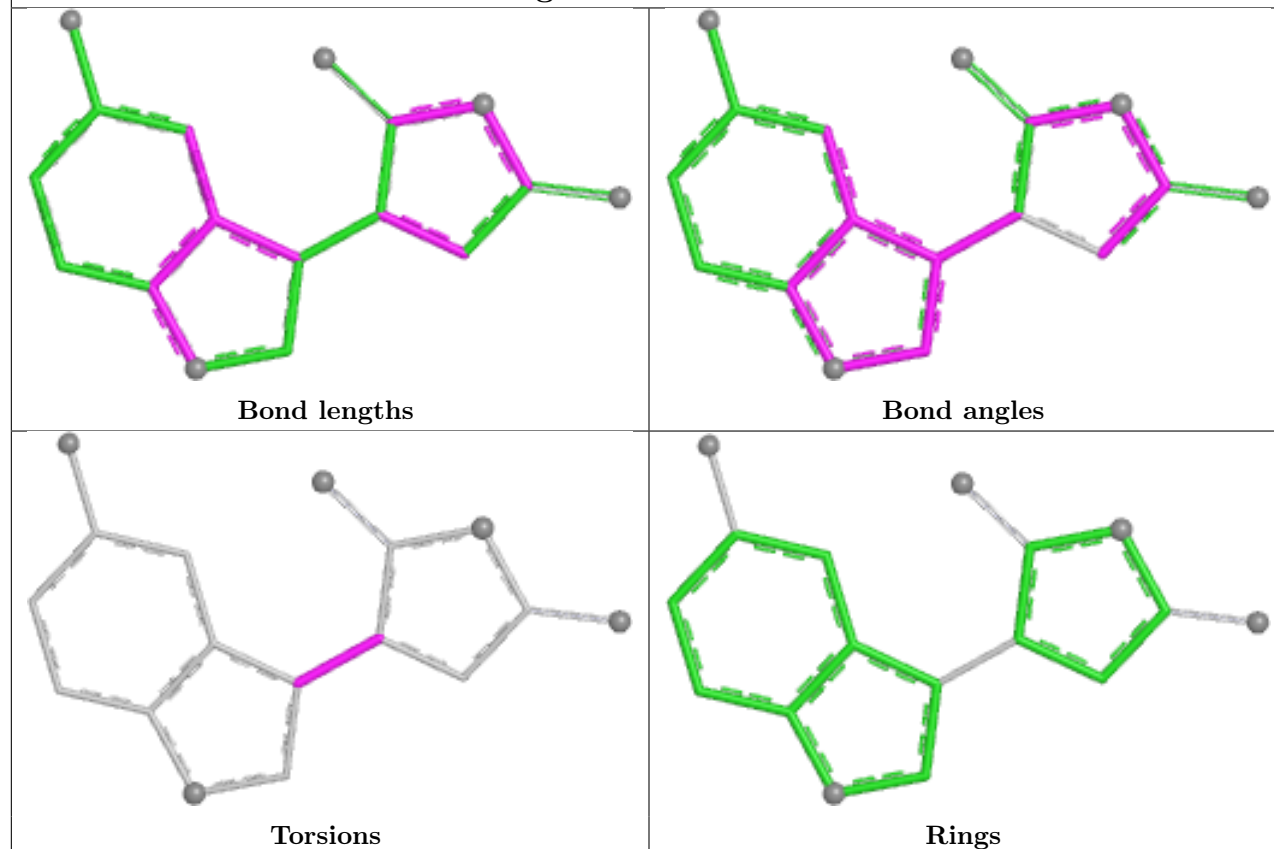
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	401	HEM	5	0
3	A	403	H7S	1	0
2	C	401	HEM	4	0
3	A	402	H7S	1	0
2	B	401	HEM	3	0
2	A	401	HEM	6	0
3	C	402	H7S	1	0
3	D	403	H7S	1	0
3	B	403	H7S	1	0
3	B	402	H7S	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths,

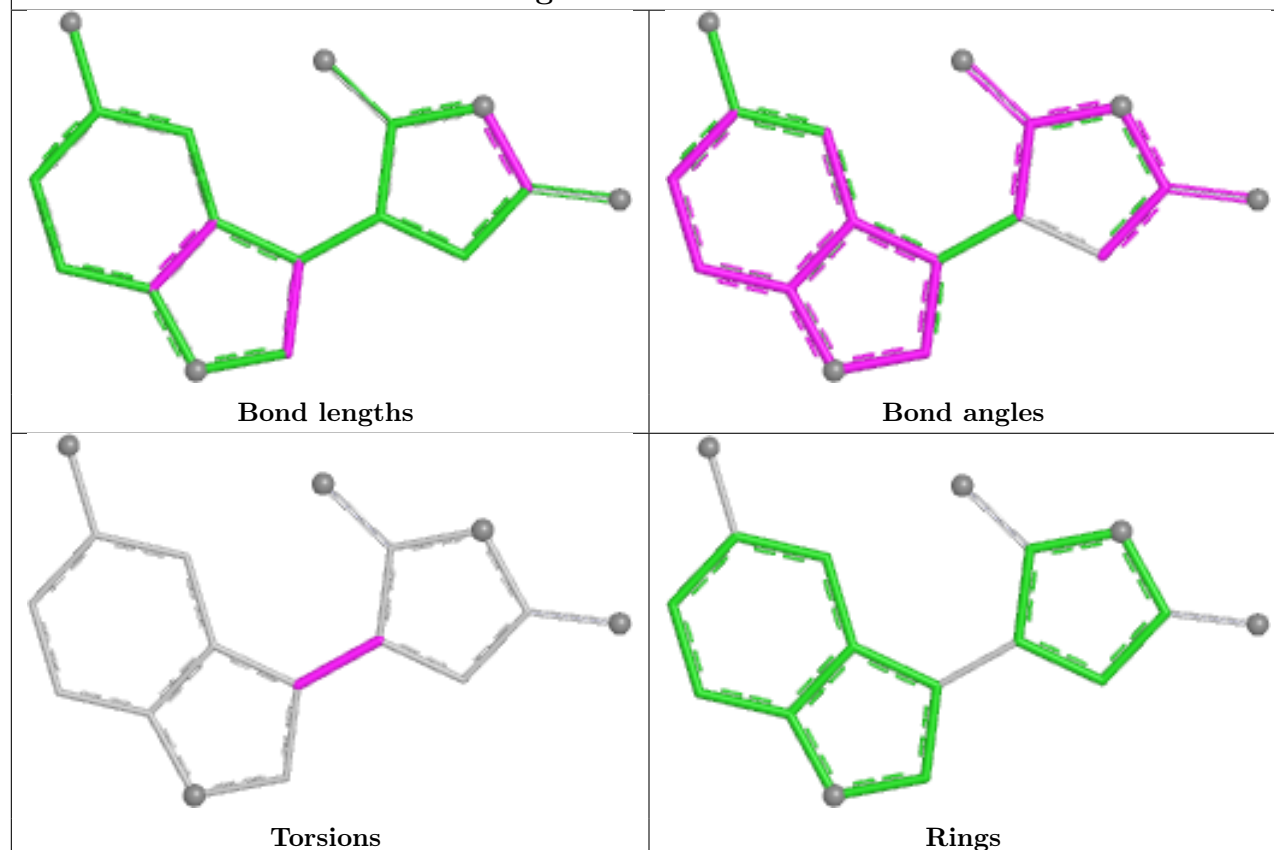
bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 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.

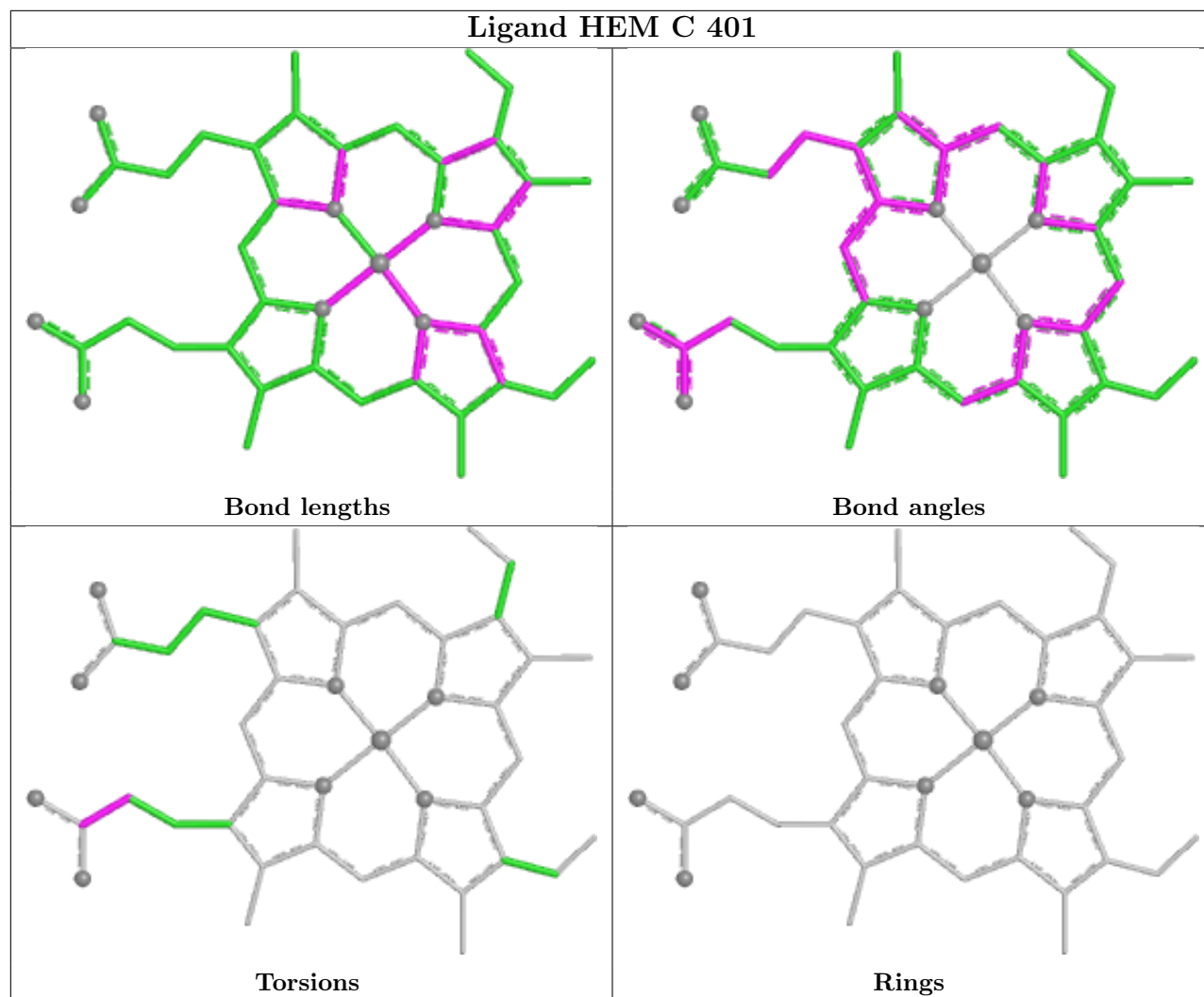


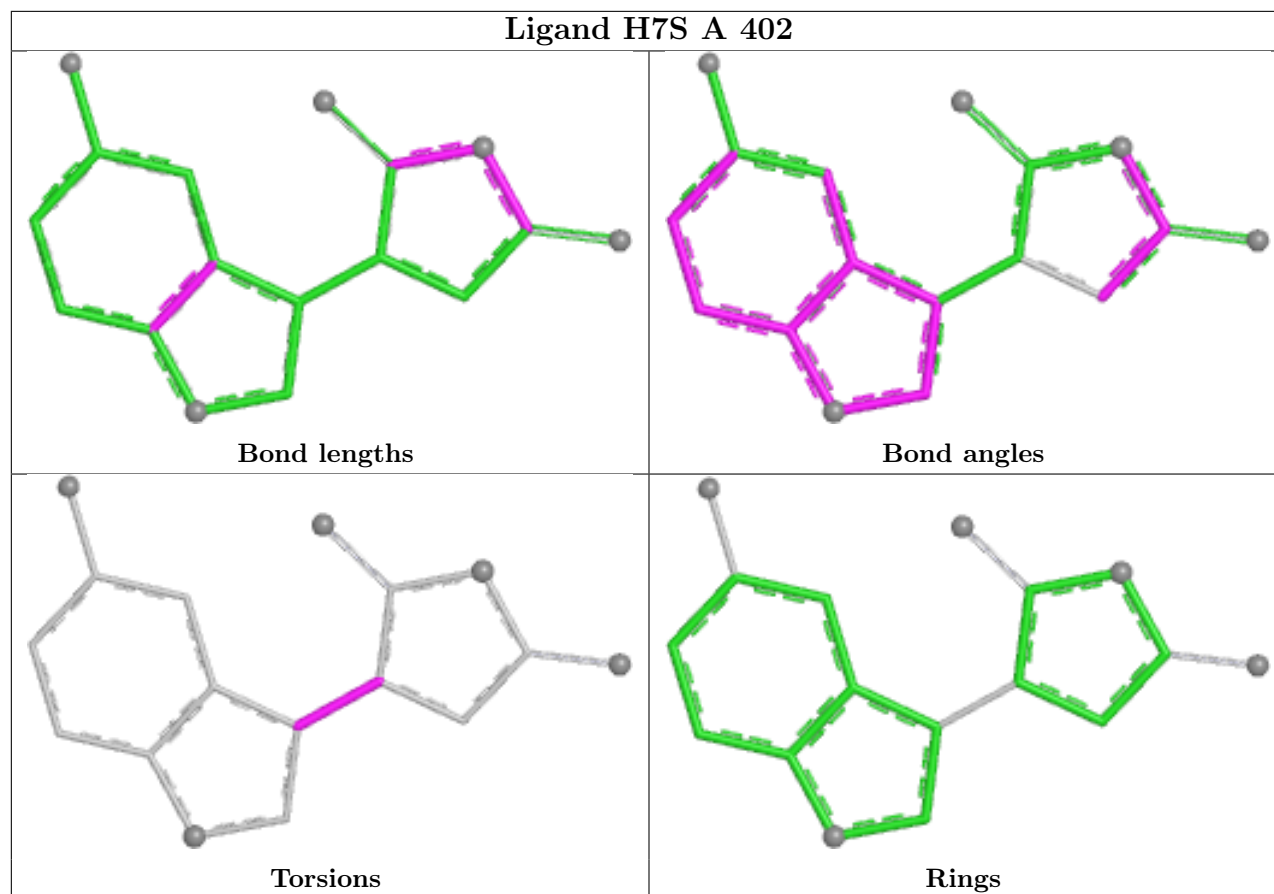
Ligand H7S A 403

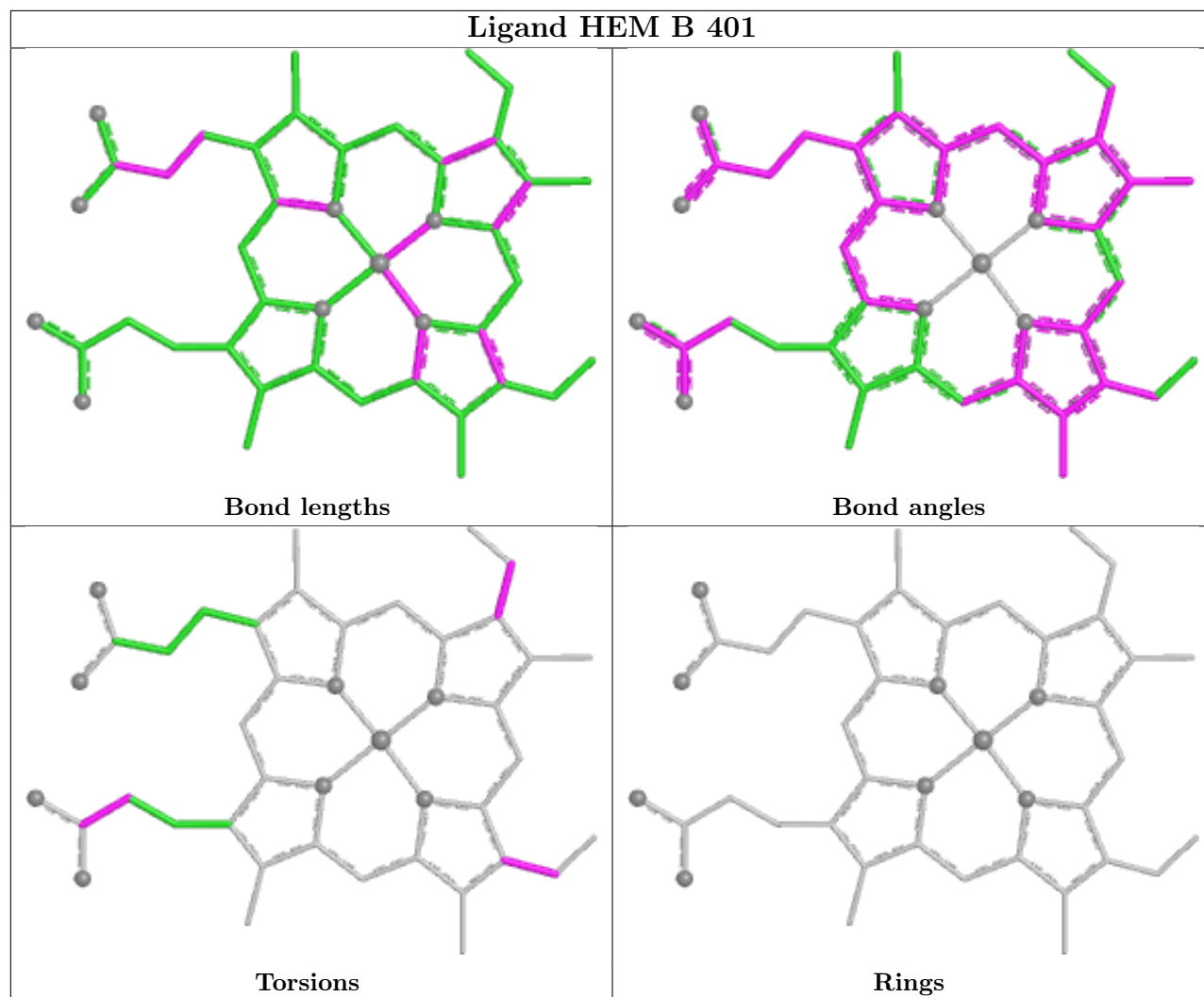


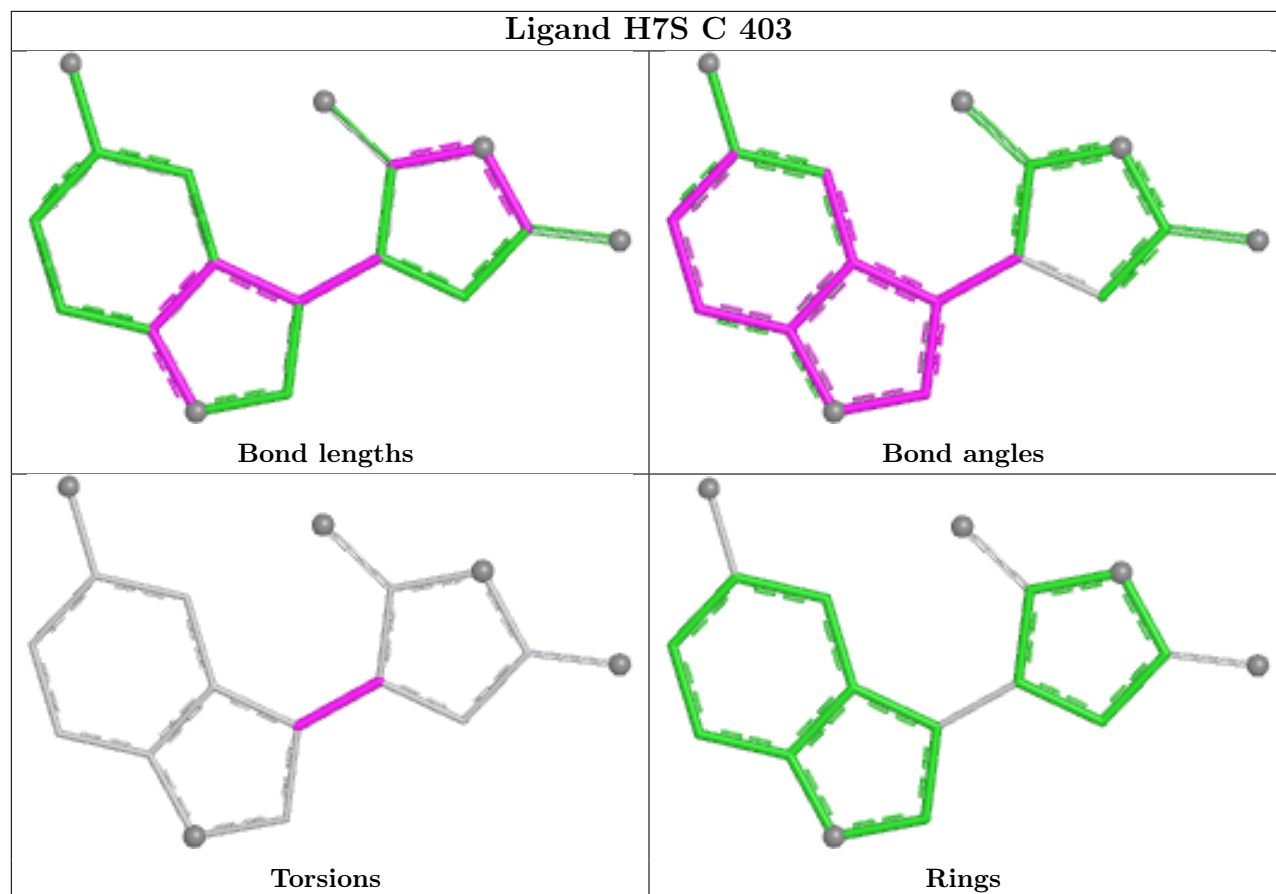
Ligand H7S D 402

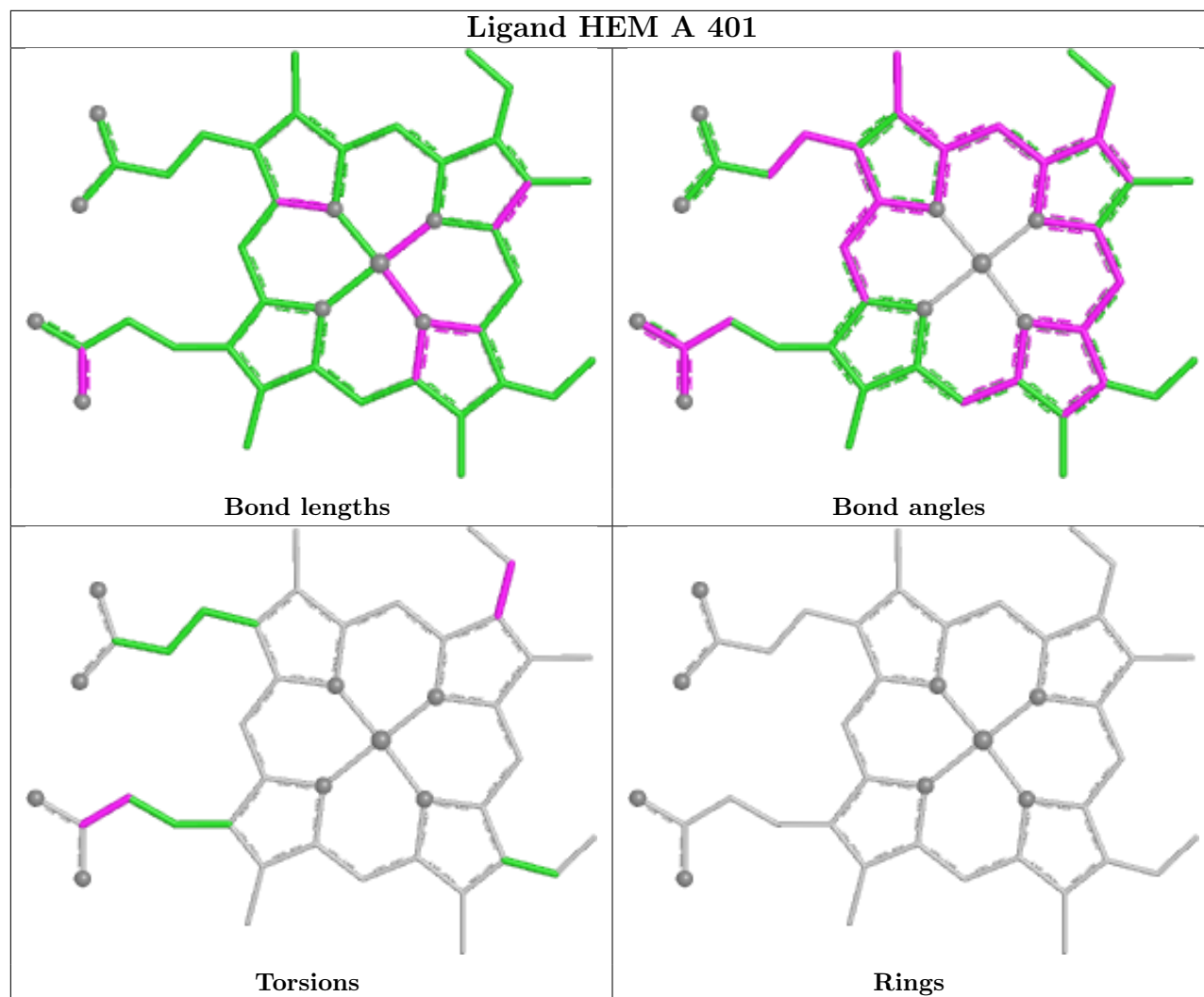




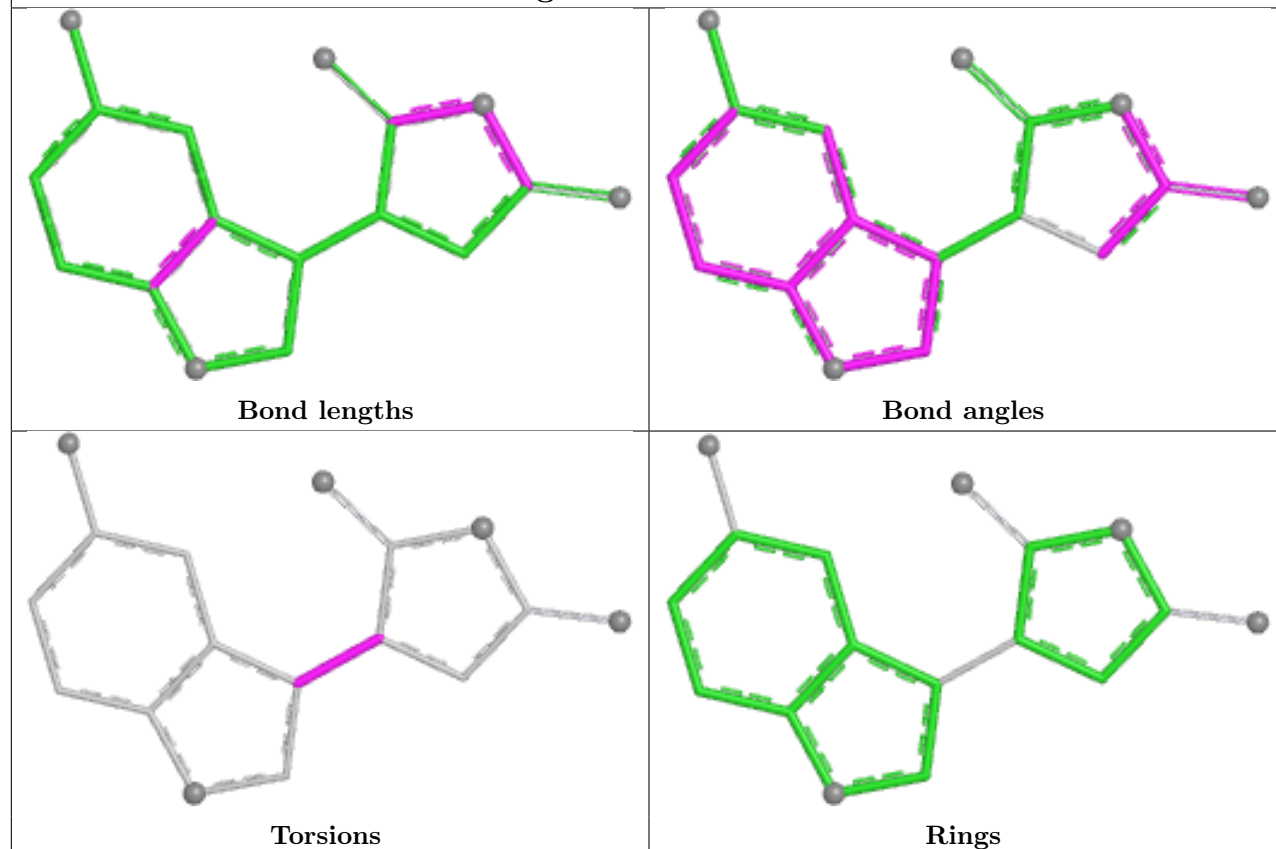




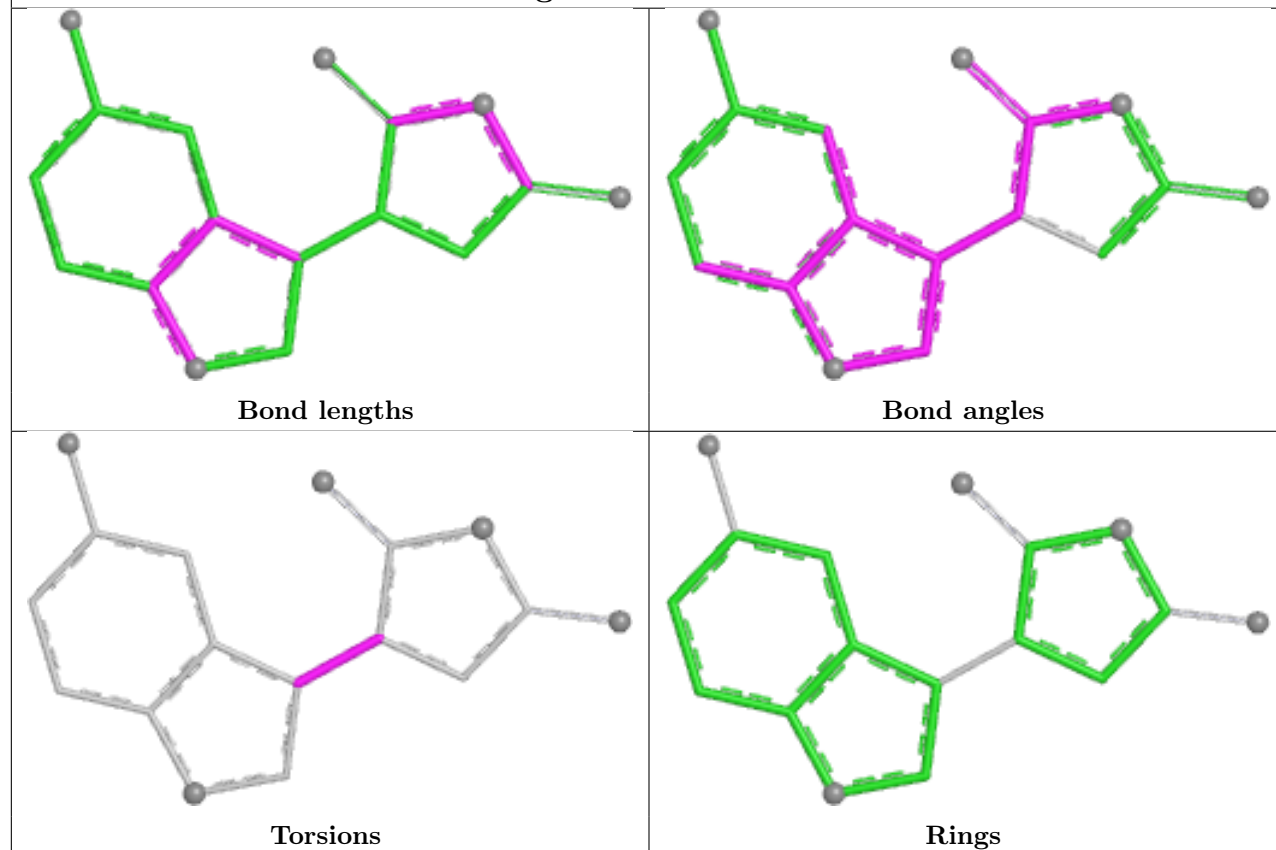




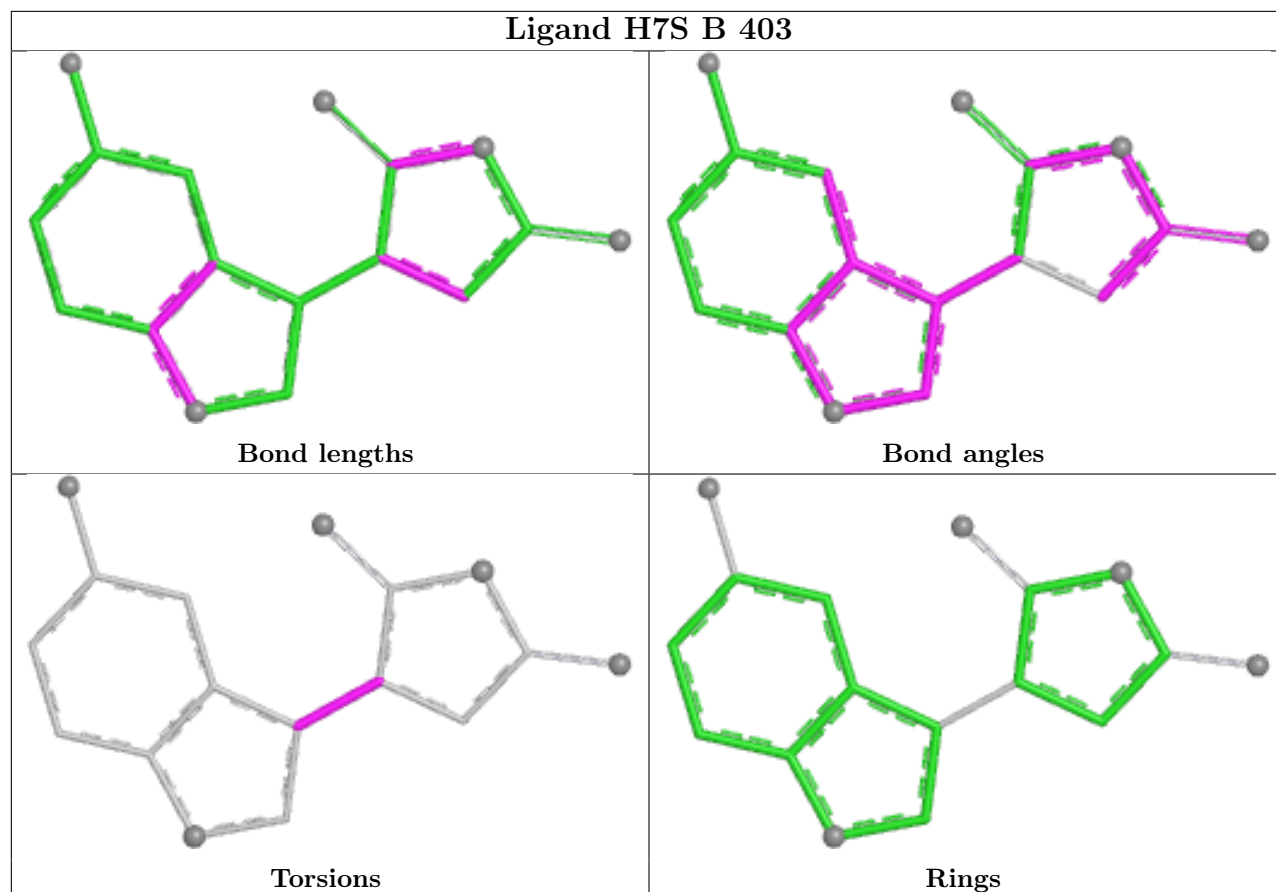
Ligand H7S C 402



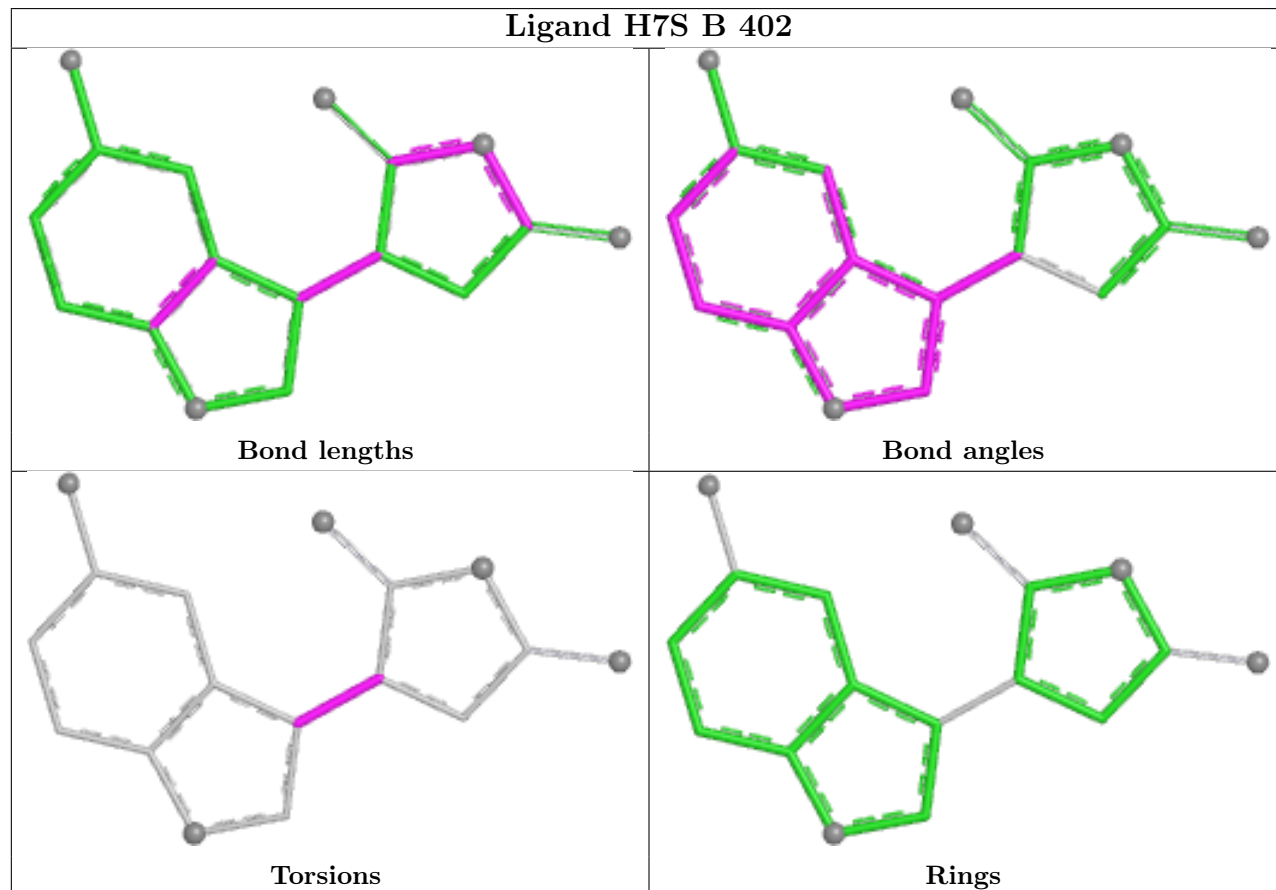
Ligand H7S D 403



Ligand H7S B 403



Ligand H7S B 402



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	351/380 (92%)	-0.03	8 (2%) 61 57	32, 50, 102, 125	0
1	B	342/380 (90%)	-0.05	9 (2%) 57 53	29, 48, 95, 140	0
1	C	332/380 (87%)	0.24	13 (3%) 43 39	33, 58, 106, 161	0
1	D	347/380 (91%)	0.05	8 (2%) 61 57	30, 52, 106, 145	0
All	All	1372/1520 (90%)	0.05	38 (2%) 55 51	29, 52, 104, 161	0

All (38) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	340	ALA	5.5
1	C	385	ILE	4.9
1	C	173	VAL	4.3
1	B	176	ASN	3.7
1	D	385	ILE	3.6
1	C	254	VAL	3.5
1	B	174	PRO	3.4
1	B	179	HIS	3.4
1	D	391	HIS	3.4
1	C	174	PRO	3.4
1	D	389	LEU	3.2
1	C	179	HIS	3.1
1	C	230	ILE	3.1
1	A	385	ILE	3.0
1	D	105	GLU	3.0
1	A	255	ALA	3.0
1	C	255	ALA	2.9
1	C	343	GLY	2.7
1	B	348	TYR	2.7
1	D	383	PRO	2.5
1	A	384	THR	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	339	LYS	2.4
1	B	173	VAL	2.4
1	B	197	GLN	2.4
1	D	175	TYR	2.3
1	A	340	ALA	2.3
1	B	170	ASN	2.3
1	D	339	LYS	2.2
1	C	345	SER	2.2
1	A	169	GLN	2.1
1	A	147	LEU	2.1
1	C	215	LEU	2.0
1	C	228	LYS	2.0
1	D	217	PRO	2.0
1	C	388	PHE	2.0
1	A	64	ASN	2.0
1	B	171	MET	2.0
1	B	376	HIS	2.0

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 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(Å ²)	Q<0.9
3	H7S	C	403	17/17	0.87	0.12	61,67,97,105	0
3	H7S	D	403	17/17	0.88	0.12	58,70,112,116	0
3	H7S	B	403	17/17	0.90	0.11	48,60,89,96	0
3	H7S	C	402	17/17	0.90	0.13	61,73,85,86	0
3	H7S	A	403	17/17	0.91	0.11	48,53,79,84	0
3	H7S	A	402	17/17	0.92	0.11	60,70,80,84	0

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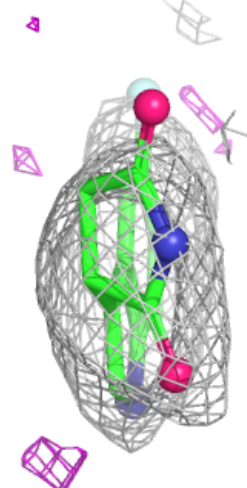
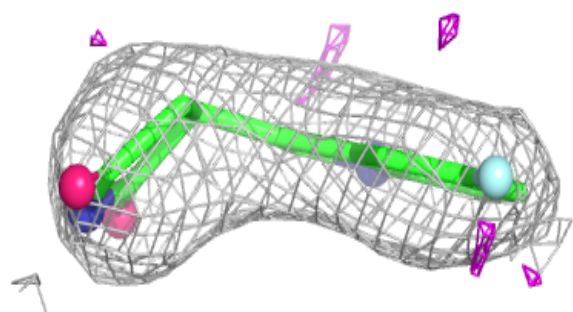
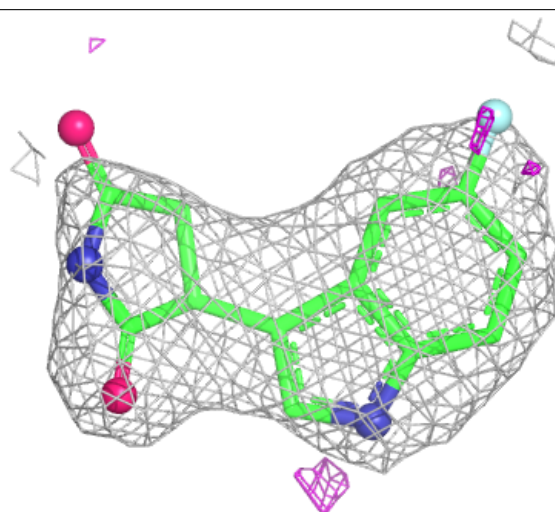
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	H7S	D	402	17/17	0.93	0.09	58,63,71,72	0
3	H7S	B	402	17/17	0.94	0.09	51,58,71,75	0
2	HEM	C	401	43/43	0.96	0.08	45,52,58,64	0
2	HEM	B	401	43/43	0.97	0.07	37,43,51,55	0
2	HEM	A	401	43/43	0.97	0.08	40,47,53,60	0
2	HEM	D	401	43/43	0.98	0.06	37,42,56,71	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.

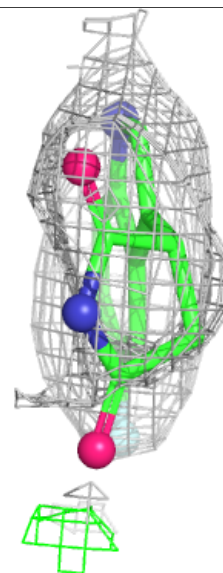
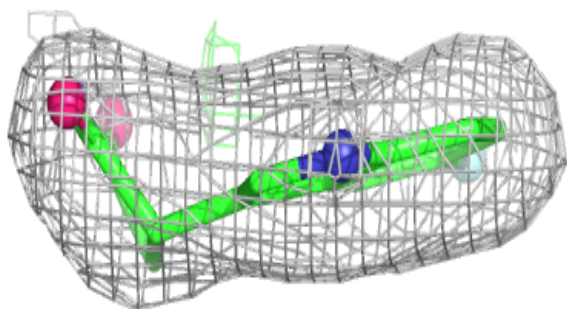
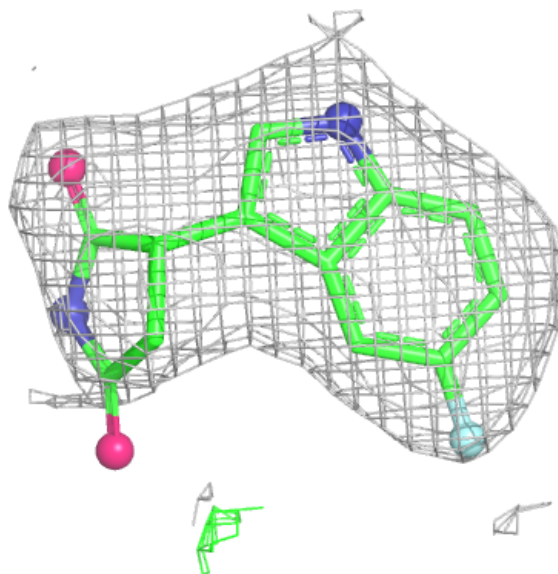
Electron density around H7S C 403:

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



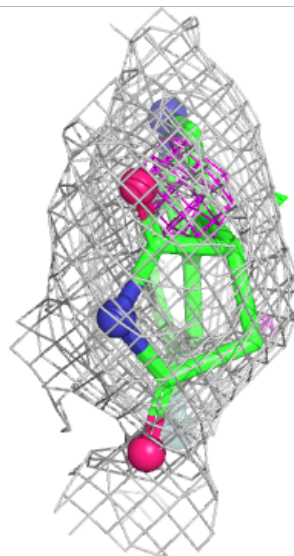
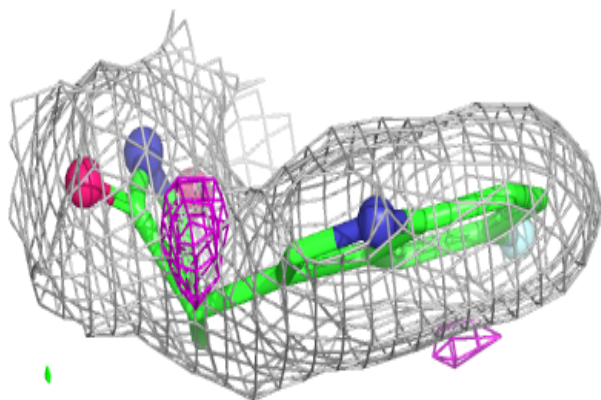
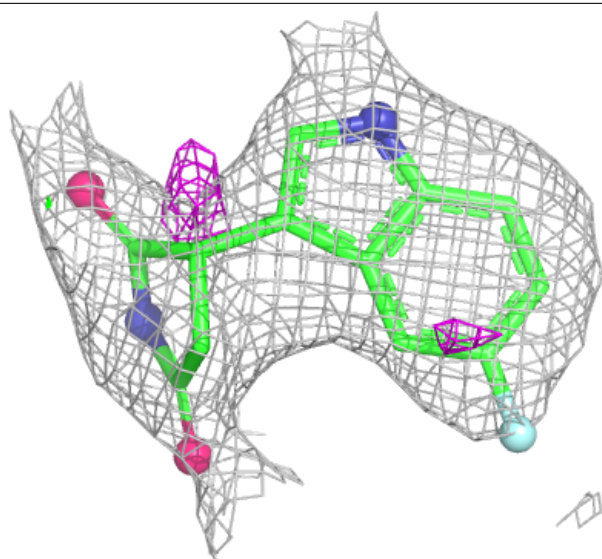
Electron density around H7S D 403:

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



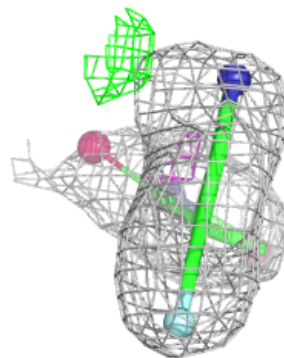
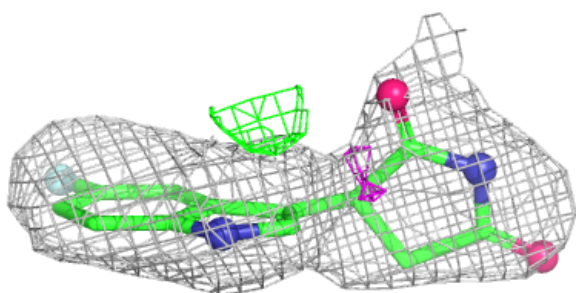
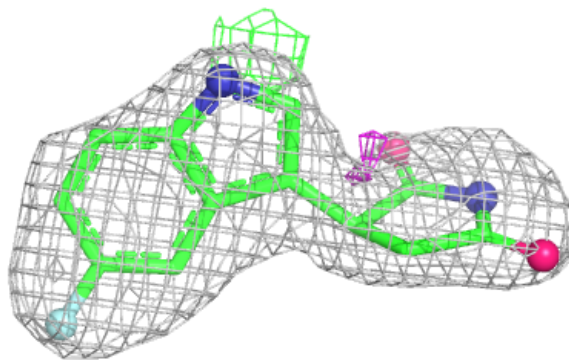
Electron density around H7S B 403:

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



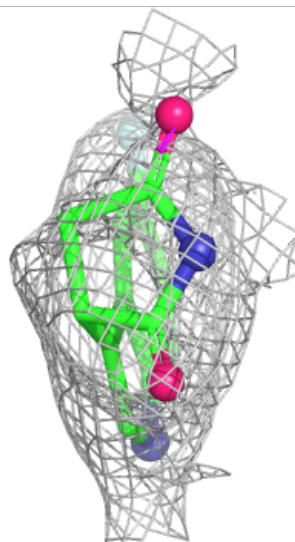
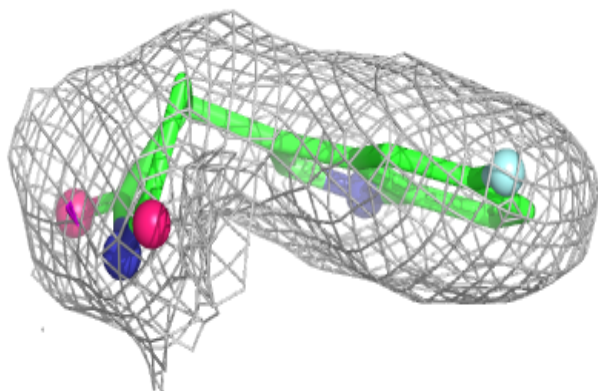
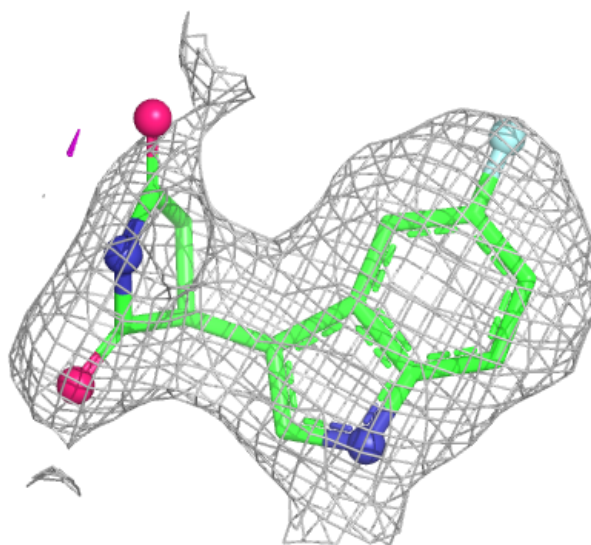
Electron density around H7S C 402:

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



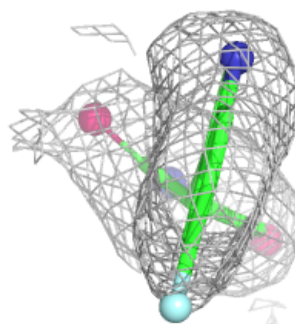
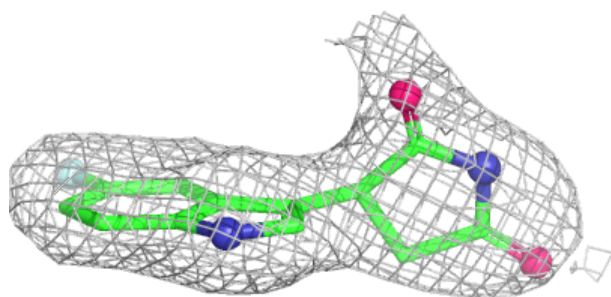
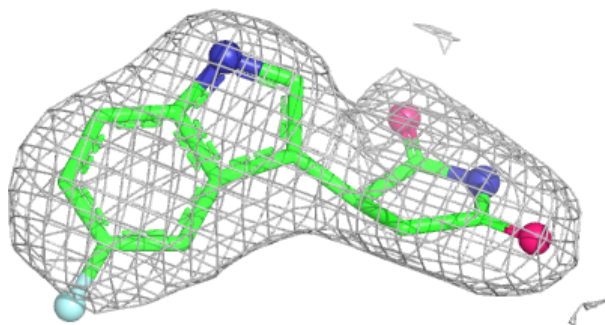
Electron density around H7S A 403:

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

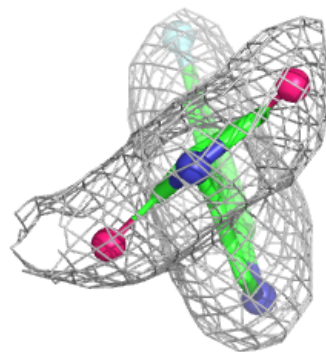
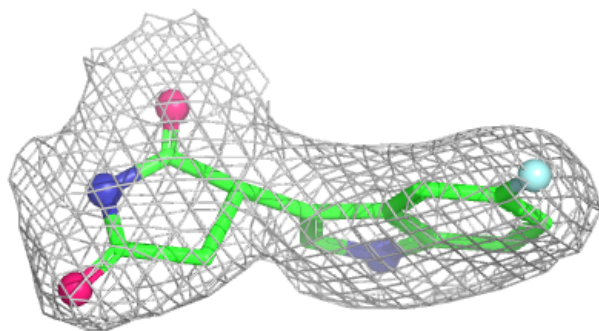
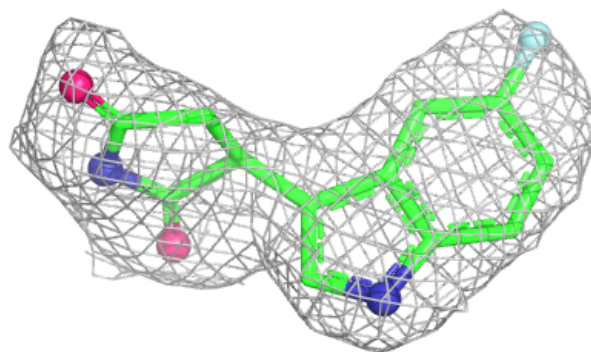


Electron density around H7S A 402:

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

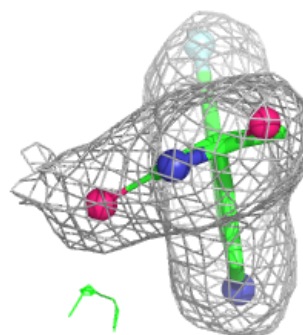
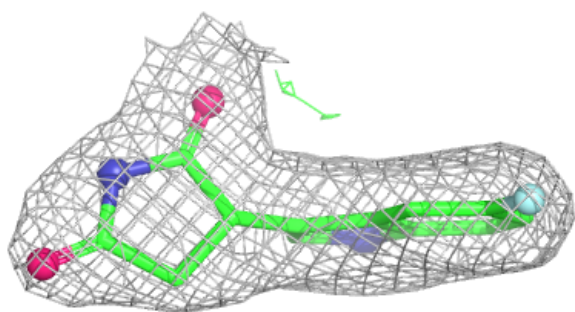
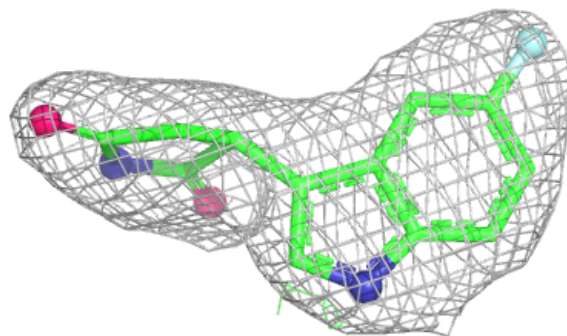
**Electron density around H7S D 402:**

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



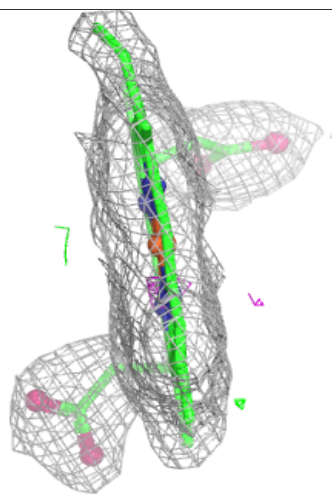
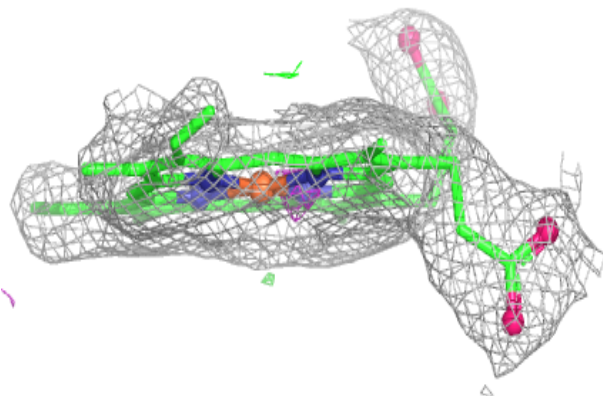
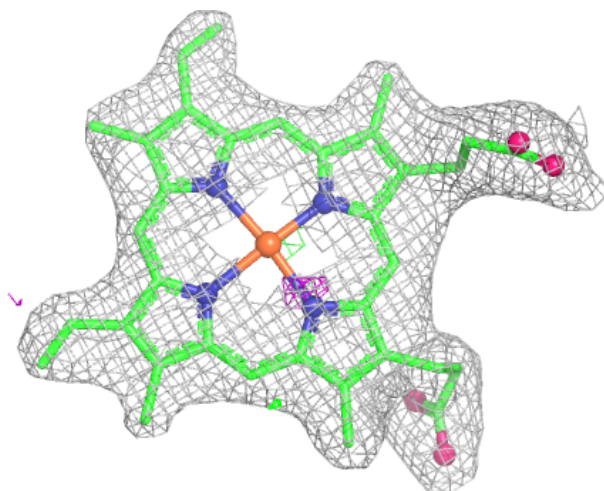
Electron density around H7S B 402:

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



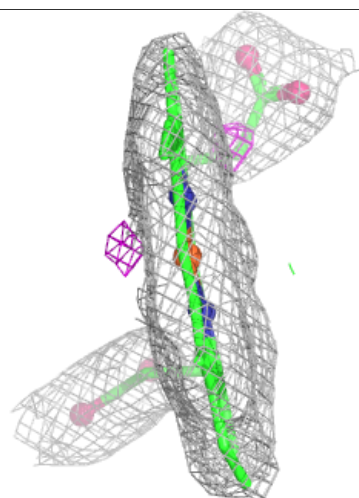
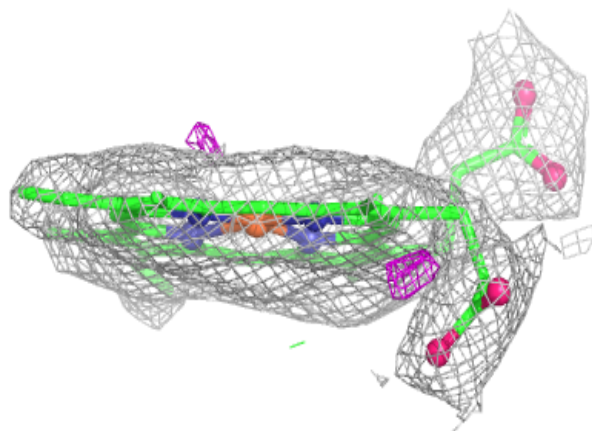
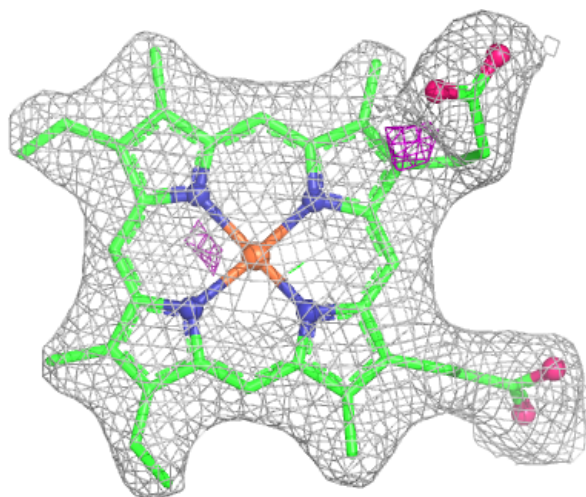
Electron density around HEM C 401:

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



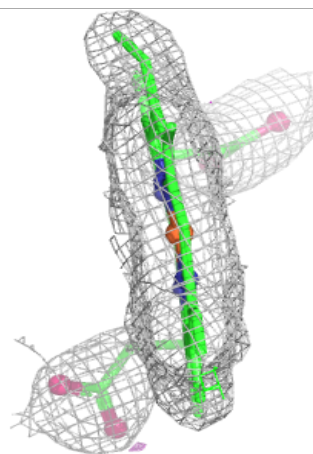
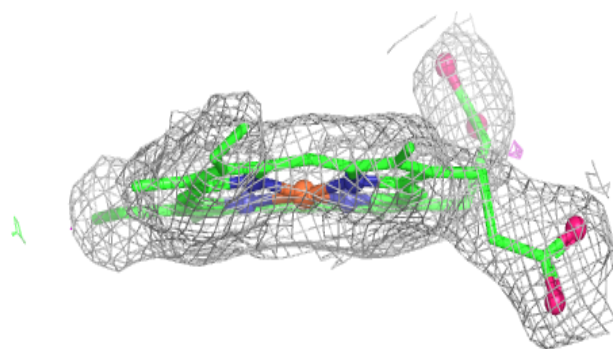
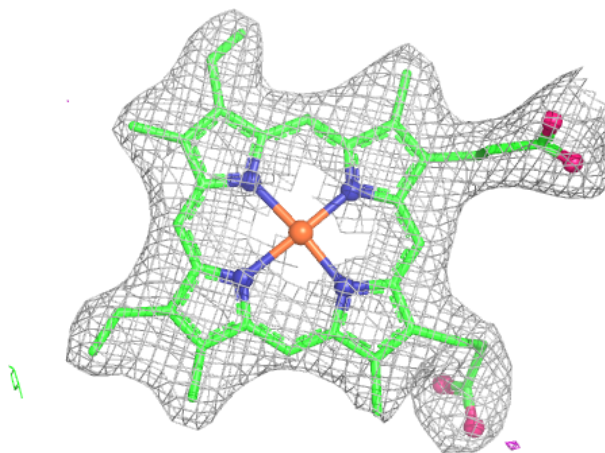
Electron density around HEM B 401:

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



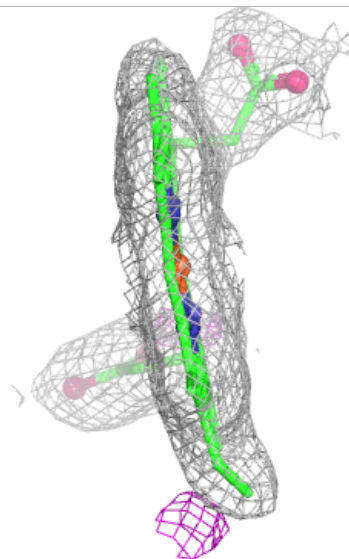
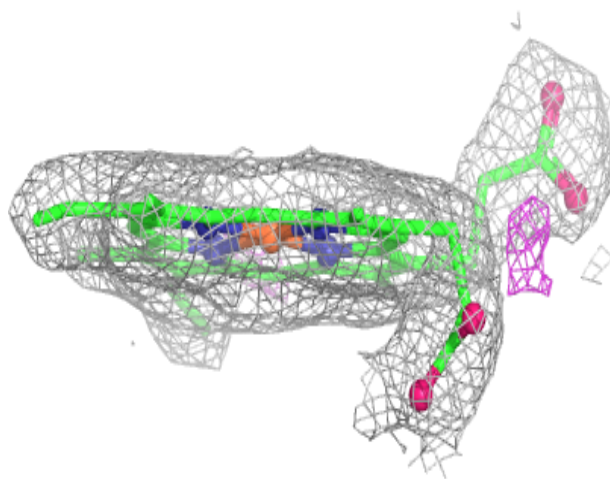
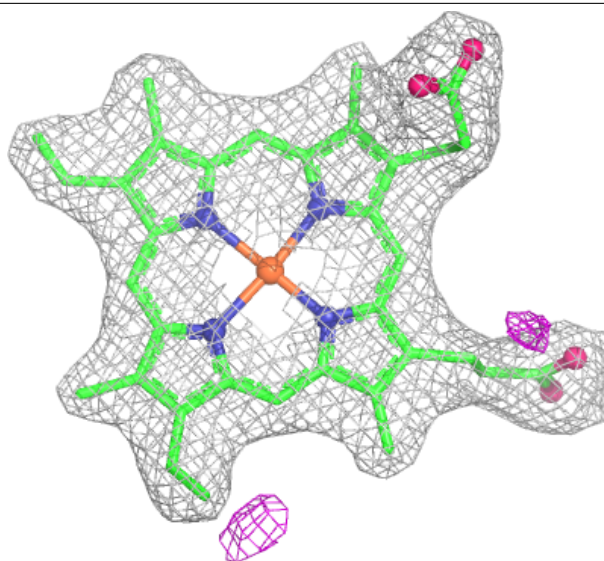
Electron density around HEM A 401:

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



Electron density around HEM D 401:

$2mF_o-DF_c$ (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.