



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 07:23 AM UTC

PDB ID : 7YJH / pdb_00007yjh
Title : Crystal structure of the P450 BM3 heme domain mutant F87V/T268I in complex with N-imidazolyl-pentanoyl-L-phenylalanine and hydroxylamine
Authors : Dong, S.; Chen, J.; Jiang, Y.; Cong, Z.; Feng, Y.
Deposited on : 2022-07-20
Resolution : 1.79 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

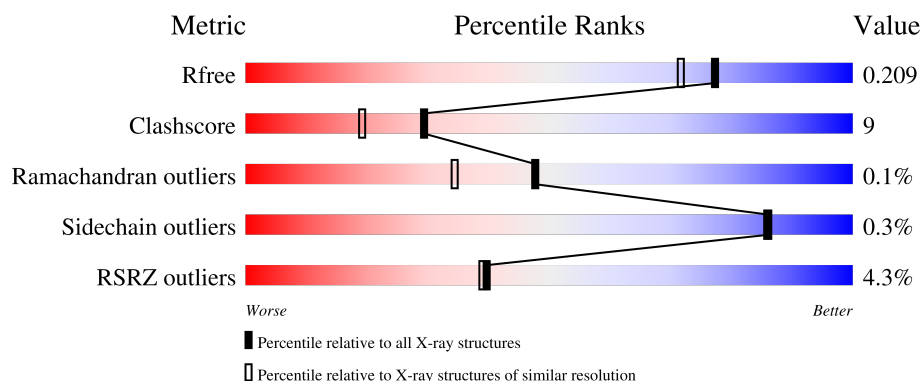
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.79 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	7662 (1.80-1.80)
Clashscore	190562	8479 (1.80-1.80)
Ramachandran outliers	187476	8391 (1.80-1.80)
Sidechain outliers	187428	8390 (1.80-1.80)
RSRZ outliers	180081	7663 (1.80-1.80)

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	466	3% 78% 19% .
1	B	466	5% 75% 21% .

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 8200 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.

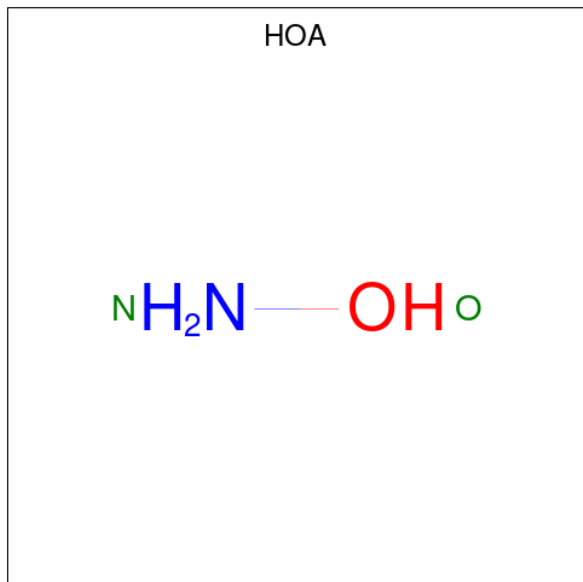
- Molecule 1 is a protein called Bifunctional cytochrome P450/NADPH-P450 reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	451	Total	C	N	O	S	0	1	0
			3625	2321	617	670	17			
1	B	451	Total	C	N	O	S	0	1	0
			3627	2321	618	671	17			

There are 24 discrepancies between the modelled and reference sequences:

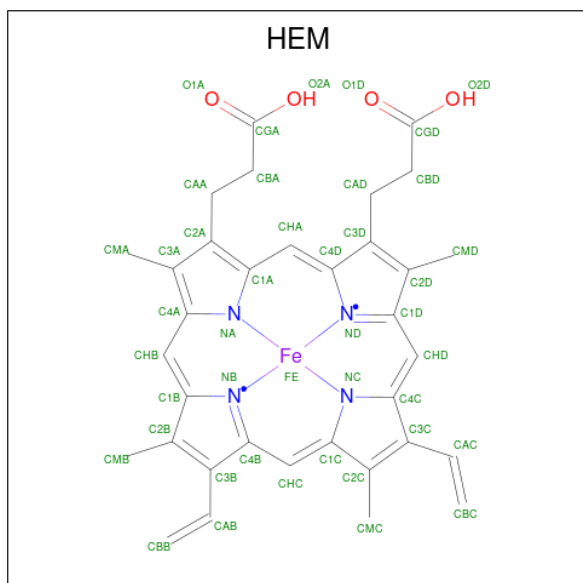
Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	MET	-	initiating methionine	UNP P14779
A	-1	GLY	-	expression tag	UNP P14779
A	87	VAL	PHE	engineered mutation	UNP P14779
A	268	ILE	THR	engineered mutation	UNP P14779
A	456	LEU	-	expression tag	UNP P14779
A	457	GLU	-	expression tag	UNP P14779
A	458	HIS	-	expression tag	UNP P14779
A	459	HIS	-	expression tag	UNP P14779
A	460	HIS	-	expression tag	UNP P14779
A	461	HIS	-	expression tag	UNP P14779
A	462	HIS	-	expression tag	UNP P14779
A	463	HIS	-	expression tag	UNP P14779
B	-2	MET	-	initiating methionine	UNP P14779
B	-1	GLY	-	expression tag	UNP P14779
B	87	VAL	PHE	engineered mutation	UNP P14779
B	268	ILE	THR	engineered mutation	UNP P14779
B	456	LEU	-	expression tag	UNP P14779
B	457	GLU	-	expression tag	UNP P14779
B	458	HIS	-	expression tag	UNP P14779
B	459	HIS	-	expression tag	UNP P14779
B	460	HIS	-	expression tag	UNP P14779
B	461	HIS	-	expression tag	UNP P14779
B	462	HIS	-	expression tag	UNP P14779
B	463	HIS	-	expression tag	UNP P14779

- Molecule 2 is HYDROXYAMINE (CCD ID: HOA) (formula: H_3NO) (labeled as "Ligand of Interest" by depositor).



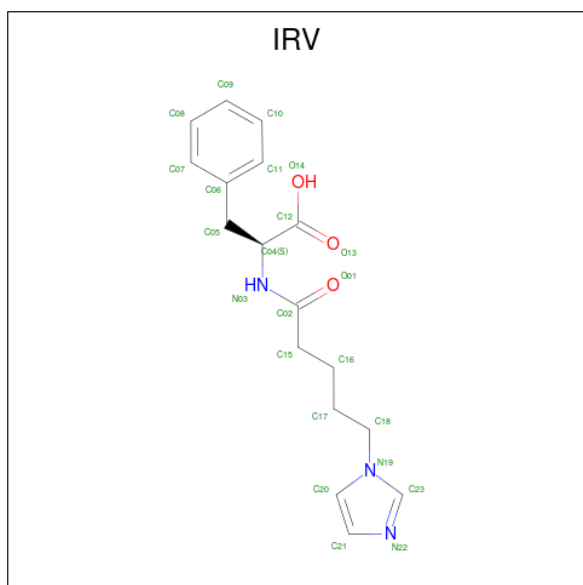
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	N	O	0	0
			2	1	1		
2	B	1	Total	N	O	0	0
			2	1	1		

- Molecule 3 is PROTOPORPHYRIN IX CONTAINING FE (CCD ID: HEM) (formula: $\text{C}_{34}\text{H}_{32}\text{FeN}_4\text{O}_4$) (labeled as "Ligand of Interest" by depositor).



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

- Molecule 4 is (2 {S})-2-(5-imidazol-1-ylpentanoylamino)-3-phenyl-propanoic acid (CCD ID: IRV) (formula: C₁₇H₂₁N₃O₃) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	1
			46	34	6	6		
4	B	1	Total	C	N	O	0	1
			46	34	6	6		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	383	Total	O	0	0
			383	383		
5	B	383	Total	O	0	0
			383	383		

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	57.27Å 144.40Å 63.41Å 90.00° 99.92° 90.00°	Depositor
Resolution (Å)	38.79 – 1.79 38.79 – 1.79	Depositor EDS
% Data completeness (in resolution range)	87.2 (38.79-1.79) 86.1 (38.79-1.79)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.66 (at 1.80Å)	Xtriage
Refinement program	PHENIX 1.13_2998	Depositor
R, R_{free}	0.167 , 0.207 0.171 , 0.209	Depositor DCC
R_{free} test set	1909 reflections (2.17%)	wwPDB-VP
Wilson B-factor (Å ²)	27.9	Xtriage
Anisotropy	0.311	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 40.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	8200	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.88% 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: HEM, IRV, HOA

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	0.46	0/3709	0.64	2/5017 (0.0%)
1	B	0.45	1/3711 (0.0%)	0.64	0/5020
All	All	0.45	1/7420 (0.0%)	0.64	2/10037 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	436	THR	CA-C	-8.16	1.49	1.53

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	436	THR	N-CA-C	-5.42	100.37	108.34
1	A	437	LEU	N-CA-C	-5.20	106.78	113.23

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	3625	0	3601	61	0
1	B	3627	0	3603	69	0
2	A	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	2	0	0	1	0
3	A	43	0	30	2	0
3	B	43	0	30	3	0
4	A	46	0	0	3	0
4	B	46	0	0	1	0
5	A	383	0	0	3	0
5	B	383	0	0	0	0
All	All	8200	0	7264	127	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 9.

All (127) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:20:LEU:O	1:A:189:GLN:HG2	1.85	0.75
1:A:148:LEU:HD21	1:A:413:VAL:HG21	1.75	0.68
1:A:104:LEU:CD1	1:B:383:SER:HB2	2.24	0.68
1:B:122:ILE:HG22	1:B:148:LEU:HD22	1.78	0.66
1:A:149:THR:HA	1:A:409:GLU:OE2	1.96	0.65
1:A:388:HIS:HA	1:A:391:LYS:HD3	1.79	0.65
1:B:367:TRP:HB2	1:B:371:VAL:HG12	1.79	0.64
1:A:22:THR:O	1:A:189:GLN:NE2	2.33	0.62
1:A:104:LEU:HD11	1:B:383:SER:HB2	1.82	0.62
1:B:81:PHE:HB3	1:B:209:ILE:HG12	1.81	0.62
1:B:163:ASN:HB2	1:B:167:ARG:HH21	1.66	0.61
1:B:379:PHE:CD2	1:B:385:ILE:HD11	2.35	0.61
1:A:120:VAL:HG11	1:A:302:VAL:HG13	1.82	0.60
1:B:214:ASP:O	1:B:218:LYS:HG3	2.02	0.60
1:B:33:ALA:CB	1:B:359:GLN:HG2	2.32	0.59
1:A:124:VAL:HG13	1:A:455:LEU:HD13	1.82	0.59
1:B:434:LYS:HB2	1:B:442:GLU:HB2	1.84	0.59
1:B:98:LYS:HD3	1:B:249:LEU:HD23	1.85	0.59
1:B:148:LEU:C	1:B:148:LEU:HD13	2.27	0.59
1:B:126:LEU:HD21	1:B:145:MET:HE1	1.85	0.58
1:B:366:ILE:HG21	1:B:389:ALA:HB1	1.84	0.58
1:A:231:ASP:OD1	1:A:236:HIS:NE2	2.35	0.58
1:A:366:ILE:HG21	1:A:389:ALA:HB1	1.84	0.58
1:A:328:ALA:O	1:A:357:ILE:HD12	2.03	0.58
1:B:298:LEU:HD22	1:B:303:PRO:HB3	1.85	0.57
1:B:149:THR:HA	1:B:409:GLU:OE2	2.06	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:98:LYS:HE2	1:B:247:GLU:HB2	1.87	0.56
1:A:81:PHE:HB3	1:A:209:ILE:HG12	1.86	0.55
1:A:47:ARG:NH1	1:A:73:GLN:HB2	2.21	0.55
1:B:328:ALA:O	1:B:357:ILE:HD12	2.06	0.54
1:A:148:LEU:CD2	1:A:413:VAL:HG21	2.39	0.53
1:B:122:ILE:HG22	1:B:148:LEU:CD2	2.39	0.53
1:A:167:ARG:NH2	5:A:617:HOH:O	2.42	0.53
1:B:286:VAL:HG11	1:B:374:PHE:HE2	1.74	0.53
1:A:419:LYS:O	1:A:451:LYS:HD2	2.09	0.52
1:B:183:GLU:O	1:B:187:LYS:HG3	2.09	0.52
1:A:74:ALA:HB2	1:A:188:LEU:HD11	1.91	0.52
1:A:104:LEU:HD13	1:B:383:SER:HB2	1.92	0.52
1:B:47:ARG:NH2	1:B:73:GLN:HB2	2.23	0.52
1:B:211:VAL:O	1:B:215:LEU:HD22	2.10	0.51
1:B:400:CYS:HA	3:B:501:HEM:CHA	2.41	0.51
1:A:329:PRO:O	1:A:358:PRO:HD3	2.11	0.51
1:A:47:ARG:NH2	4:A:503[B]:IRV:O14	2.40	0.51
1:B:215:LEU:O	1:B:219:ILE:HG12	2.11	0.51
1:A:429:TYR:CE2	1:A:431:LEU:HA	2.46	0.51
1:A:434:LYS:HB2	1:A:442:GLU:HB2	1.93	0.50
1:B:198:TYR:O	1:B:202:LYS:HG3	2.12	0.50
1:B:409:GLU:O	1:B:413:VAL:HG12	2.12	0.50
2:B:502:HOA:O	4:B:503[B]:IRV:C21	2.59	0.50
1:A:298:LEU:HD22	1:A:303:PRO:HB3	1.92	0.49
1:A:59:LYS:HE3	1:B:296:ARG:HD3	1.94	0.49
1:A:236:HIS:O	1:A:240:GLY:N	2.46	0.49
1:B:122:ILE:CG2	1:B:148:LEU:HD22	2.42	0.49
1:A:22:THR:H	1:A:189:GLN:NE2	2.11	0.48
1:A:211:VAL:O	1:A:215:LEU:HG	2.13	0.48
1:B:278:TYR:HA	1:B:444:PHE:CZ	2.49	0.48
1:B:24:LYS:HE2	1:B:433:ILE:O	2.14	0.48
1:A:272:LEU:HD13	1:A:322:LEU:HG	1.96	0.48
1:A:317:VAL:HG13	1:A:374:PHE:HZ	1.79	0.47
1:B:253:ASN:O	1:B:257:GLN:HG2	2.15	0.47
1:B:69:LYS:NZ	1:B:87[B]:VAL:O	2.48	0.47
1:B:393:PHE:HB3	1:B:400:CYS:HB3	1.97	0.47
1:B:14:LEU:HB3	1:B:18:PRO:HD3	1.98	0.46
1:B:379:PHE:CD2	1:B:385:ILE:CD1	2.99	0.46
1:B:112:MET:SD	1:B:405:PHE:HA	2.55	0.46
1:B:272:LEU:HD13	1:B:322:LEU:HG	1.96	0.46
1:B:234:LEU:HD13	1:B:258:ILE:HD11	1.96	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:124:VAL:HG13	1:B:455:LEU:HD13	1.97	0.46
1:B:449:LYS:HD3	1:B:449:LYS:HA	1.59	0.46
1:A:320:GLU:HG3	1:A:374:PHE:CD1	2.51	0.45
1:B:110:GLN:OE1	1:B:113:LYS:HE3	2.15	0.45
1:A:271:GLY:HA3	1:A:327:THR:HG21	1.98	0.45
1:A:366:ILE:HD12	1:A:386:PRO:HG2	1.98	0.45
1:A:311:LEU:HB3	1:A:314:VAL:HB	1.98	0.45
1:B:241:LYS:NZ	1:B:248:PRO:HB3	2.31	0.45
1:B:429:TYR:CE2	1:B:431:LEU:HA	2.51	0.45
1:A:278:TYR:HA	1:A:444:PHE:CZ	2.51	0.45
1:B:157:GLY:HA2	1:B:233:LEU:HD12	1.98	0.45
1:B:359:GLN:OE1	1:B:359:GLN:HA	2.16	0.45
1:A:223:ARG:NH2	1:A:232:ASP:OD2	2.42	0.45
1:B:141:VAL:O	1:B:145:MET:HG2	2.17	0.45
1:A:393:PHE:HB3	1:A:400:CYS:HB3	1.99	0.44
1:B:326:PRO:HG3	1:B:357:ILE:HG22	2.00	0.44
1:A:106:SER:HB3	1:A:233:LEU:HD23	2.00	0.44
1:A:24:LYS:HE2	1:A:433:ILE:O	2.18	0.44
1:B:130:TRP:CZ2	1:B:139:ILE:HG21	2.53	0.44
1:B:39:ILE:HA	1:B:51:TYR:O	2.18	0.44
1:A:400:CYS:HA	3:A:502:HEM:CHA	2.48	0.43
1:A:319:ASN:ND2	5:A:635:HOH:O	2.50	0.43
1:B:33:ALA:HB3	1:B:359:GLN:HG2	2.00	0.43
1:A:51:TYR:OH	4:A:503[A]:IRV:O14	2.31	0.43
1:B:33:ALA:HB1	1:B:359:GLN:HG2	1.99	0.43
1:A:75:LEU:HD23	1:A:75:LEU:HA	1.85	0.43
1:A:130:TRP:CZ2	1:A:139:ILE:HG21	2.53	0.42
1:B:320:GLU:HG3	1:B:374:PHE:CD1	2.54	0.42
1:B:16:ASN:HB3	1:B:19:LEU:HD12	2.00	0.42
1:B:106:SER:HB3	1:B:233:LEU:HD23	2.01	0.42
1:A:280:LEU:HD22	1:A:287:LEU:HA	2.01	0.42
1:B:68:ASP:HB3	1:B:334:TYR:CZ	2.54	0.42
1:B:153:ILE:HD12	1:B:153:ILE:HA	1.90	0.42
1:A:190:ARG:HG2	1:A:198:TYR:CZ	2.55	0.42
1:B:400:CYS:HA	3:B:501:HEM:C4D	2.54	0.42
1:A:53:SER:O	1:A:360:LEU:HA	2.20	0.42
1:A:407:LEU:HD23	1:A:407:LEU:HA	1.84	0.42
1:A:112:MET:SD	1:A:405:PHE:HA	2.59	0.42
1:B:280:LEU:HD22	1:B:287:LEU:HA	2.02	0.42
1:B:391:LYS:HE2	1:B:395:ASN:HB2	2.01	0.42
1:A:271:GLY:C	1:A:327:THR:HG21	2.45	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:400:CYS:HA	3:A:502:HEM:C4D	2.55	0.42
1:A:323:ARG:HG2	1:A:361:HIS:HB3	2.01	0.41
1:B:150:LEU:HD21	1:B:174:ILE:HG13	2.01	0.41
1:A:331:PHE:HD2	1:A:357:ILE:HD11	1.84	0.41
1:A:452:LYS:HE3	1:A:452:LYS:HB3	1.89	0.41
1:B:153:ILE:O	1:B:157:GLY:N	2.44	0.41
1:A:20:LEU:HG	1:A:42:PHE:CZ	2.56	0.41
1:B:366:ILE:HG21	1:B:389:ALA:CB	2.48	0.41
1:A:286:VAL:HG11	1:A:374:PHE:CE2	2.56	0.41
1:A:106:SER:HB3	1:A:233:LEU:CD2	2.51	0.41
1:B:130:TRP:HB2	1:B:421:PHE:CZ	2.56	0.41
1:B:267:GLU:HA	1:B:270:SER:OG	2.20	0.41
1:A:142:PRO:HA	1:A:274:SER:OG	2.21	0.41
1:A:87[A]:VAL:HG21	4:A:503[A]:IRV:C21	2.51	0.40
1:A:440:LYS:HB3	5:A:719:HOH:O	2.20	0.40
1:B:20:LEU:HG	1:B:42:PHE:CZ	2.56	0.40
1:A:39:ILE:HA	1:A:51:TYR:O	2.21	0.40
1:B:293:GLU:CD	1:B:313:TYR:H	2.29	0.40
1:B:400:CYS:HB2	3:B:501:HEM:NA	2.36	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	450/466 (97%)	431 (96%)	19 (4%)	0	100	100
1	B	450/466 (97%)	435 (97%)	14 (3%)	1 (0%)	43	31
All	All	900/932 (97%)	866 (96%)	33 (4%)	1 (0%)	48	34

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	266	HIS

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	391/409 (96%)	390 (100%)	1 (0%)	86	86
1	B	392/409 (96%)	391 (100%)	1 (0%)	86	86
All	All	783/818 (96%)	781 (100%)	2 (0%)	86	86

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	371	VAL
1	B	359	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	73	GLN
1	A	189	GLN
1	A	204	GLN
1	A	229	GLN
1	A	239	ASN
1	B	201	ASN
1	B	204	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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](#)

8 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	HOA	B	502	3	0,1,1	-	-	-		
3	HEM	B	501	1,2	50,50,50	1.33	5 (10%)	67,82,82	1.19	6 (8%)
4	IRV	A	503[B]	-	24,24,24	1.44	1 (4%)	30,30,30	1.09	2 (6%)
4	IRV	A	503[A]	-	24,24,24	1.51	2 (8%)	30,30,30	0.95	1 (3%)
2	HOA	A	501	3	0,1,1	-	-	-		
4	IRV	B	503[B]	-	24,24,24	1.45	2 (8%)	30,30,30	1.02	2 (6%)
4	IRV	B	503[A]	-	24,24,24	1.46	2 (8%)	30,30,30	0.97	1 (3%)
3	HEM	A	502	1,2	50,50,50	1.35	6 (12%)	67,82,82	1.21	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
3	HEM	B	501	1,2	-	2/14/54/54	-
4	IRV	A	503[B]	-	-	8/19/19/19	0/2/2/2
4	IRV	A	503[A]	-	-	1/19/19/19	0/2/2/2
4	IRV	B	503[B]	-	-	2/19/19/19	0/2/2/2
4	IRV	B	503[A]	-	-	5/19/19/19	0/2/2/2
3	HEM	A	502	1,2	-	2/14/54/54	-

All (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	503[A]	IRV	C02-N03	5.86	1.46	1.34
4	A	503[B]	IRV	C02-N03	5.48	1.45	1.34
4	B	503[B]	IRV	C02-N03	5.46	1.45	1.34
4	B	503[A]	IRV	C02-N03	5.43	1.45	1.34
3	A	502	HEM	CAC-C3C	2.99	1.55	1.47
3	A	502	HEM	CAB-C3B	2.85	1.55	1.47
3	B	501	HEM	CAB-C3B	2.77	1.54	1.47
3	A	502	HEM	FE-NC	2.51	2.03	1.95
3	B	501	HEM	CAC-C3C	2.41	1.53	1.47
3	B	501	HEM	FE-NA	2.41	2.03	1.95
3	A	502	HEM	CMC-C2C	2.40	1.55	1.50
3	A	502	HEM	CMB-C2B	2.40	1.55	1.50
4	B	503[A]	IRV	O01-C02	-2.34	1.18	1.23
4	B	503[B]	IRV	O01-C02	-2.33	1.18	1.23
3	B	501	HEM	C2A-C3A	-2.29	1.32	1.38
4	A	503[A]	IRV	O01-C02	-2.17	1.19	1.23
3	A	502	HEM	FE-NB	2.13	2.01	1.94
3	B	501	HEM	CMC-C2C	2.03	1.54	1.50

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	501	HEM	CBA-CAA-C2A	-3.01	104.21	112.53
3	A	502	HEM	CBA-CAA-C2A	-2.91	104.48	112.53
4	A	503[B]	IRV	C12-C04-N03	-2.90	103.83	110.57
3	A	502	HEM	CAD-CBD-CGD	-2.74	106.40	113.67
3	B	501	HEM	CHA-C4D-ND	2.63	127.62	124.37
3	A	502	HEM	CBD-CAD-C3D	-2.59	105.37	112.53
3	B	501	HEM	CAD-CBD-CGD	-2.57	106.86	113.67
3	B	501	HEM	CHB-C4A-NA	2.56	128.51	123.86
3	B	501	HEM	CBD-CAD-C3D	-2.46	105.72	112.53
4	B	503[B]	IRV	C15-C02-N03	2.36	120.02	115.86
4	A	503[B]	IRV	O14-C12-C04	2.27	121.21	113.51
4	A	503[A]	IRV	C15-C02-N03	2.22	119.77	115.86
3	B	501	HEM	C2A-C1A-NA	-2.20	107.71	110.15
3	A	502	HEM	CHD-C1D-ND	2.15	126.74	124.42
4	B	503[A]	IRV	C15-C02-N03	2.15	119.65	115.86
4	B	503[B]	IRV	O14-C12-C04	2.11	120.64	113.51
3	A	502	HEM	C2A-C1A-NA	-2.05	107.88	110.15

There are no chirality outliers.

All (20) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	503[A]	IRV	C02-C15-C16-C17
4	B	503[B]	IRV	C16-C17-C18-N19
4	B	503[A]	IRV	N03-C02-C15-C16
4	B	503[A]	IRV	O01-C02-C15-C16
4	B	503[A]	IRV	C16-C17-C18-N19
4	A	503[B]	IRV	C02-C15-C16-C17
4	B	503[A]	IRV	C17-C18-N19-C20
4	A	503[B]	IRV	N03-C02-C15-C16
4	A	503[B]	IRV	O01-C02-C15-C16
4	A	503[B]	IRV	C12-C04-N03-C02
4	B	503[A]	IRV	C17-C18-N19-C23
4	A	503[B]	IRV	C15-C16-C17-C18
3	B	501	HEM	CAD-CBD-CGD-O1D
4	A	503[B]	IRV	C17-C18-N19-C23
4	A	503[B]	IRV	C17-C18-N19-C20
3	B	501	HEM	CAD-CBD-CGD-O2D
4	A	503[B]	IRV	C16-C17-C18-N19
3	A	502	HEM	CAD-CBD-CGD-O2D
3	A	502	HEM	CAD-CBD-CGD-O1D
4	B	503[B]	IRV	O01-C02-C15-C16

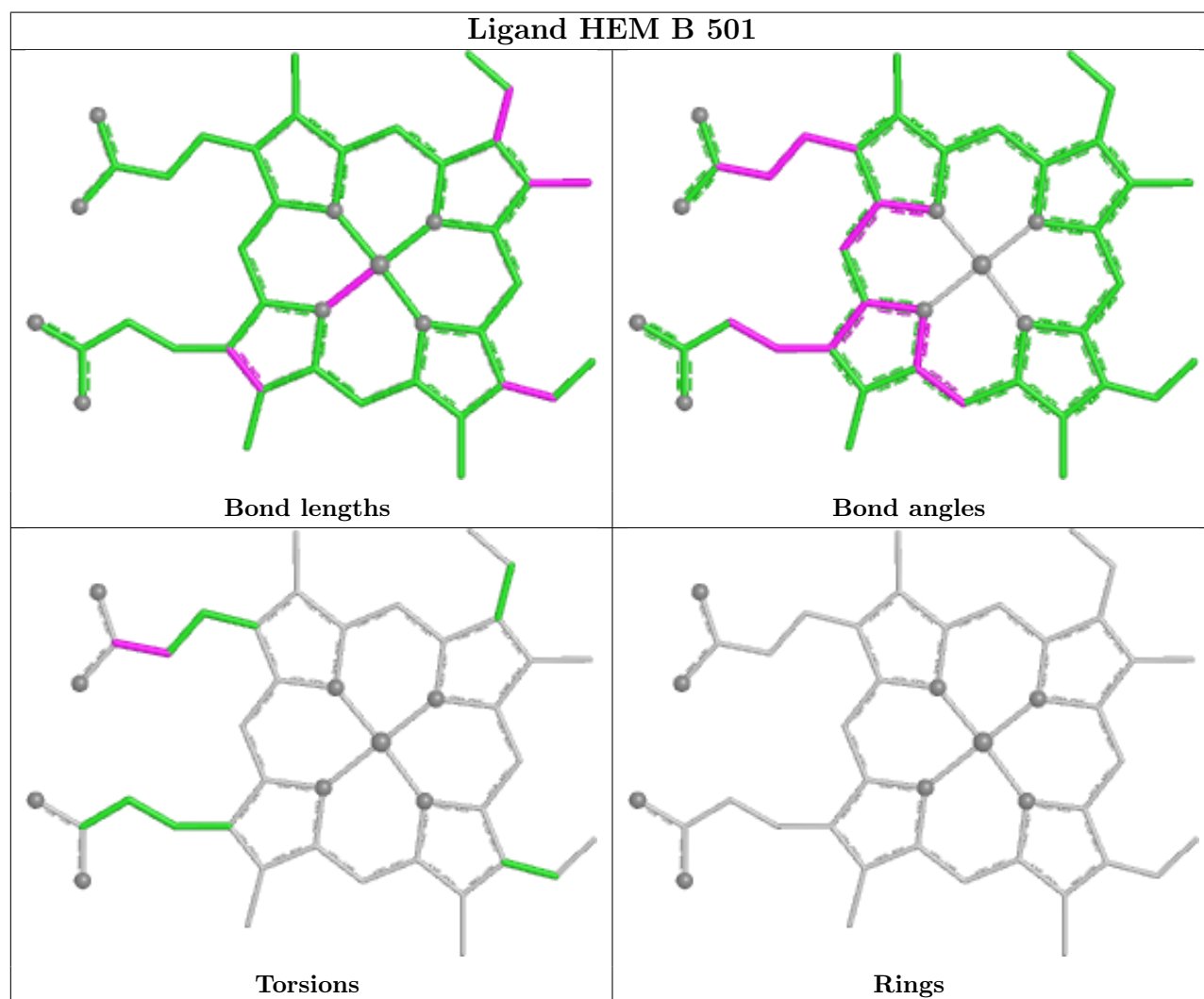
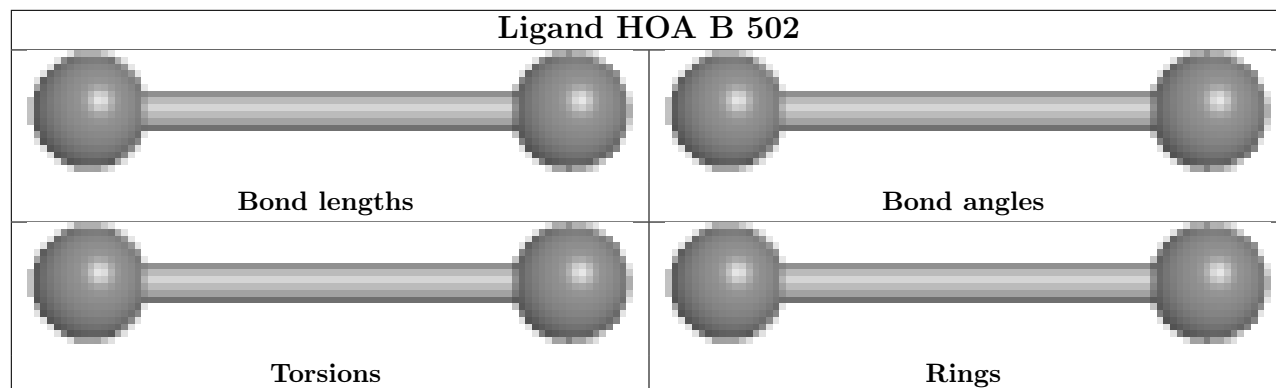
There are no ring outliers.

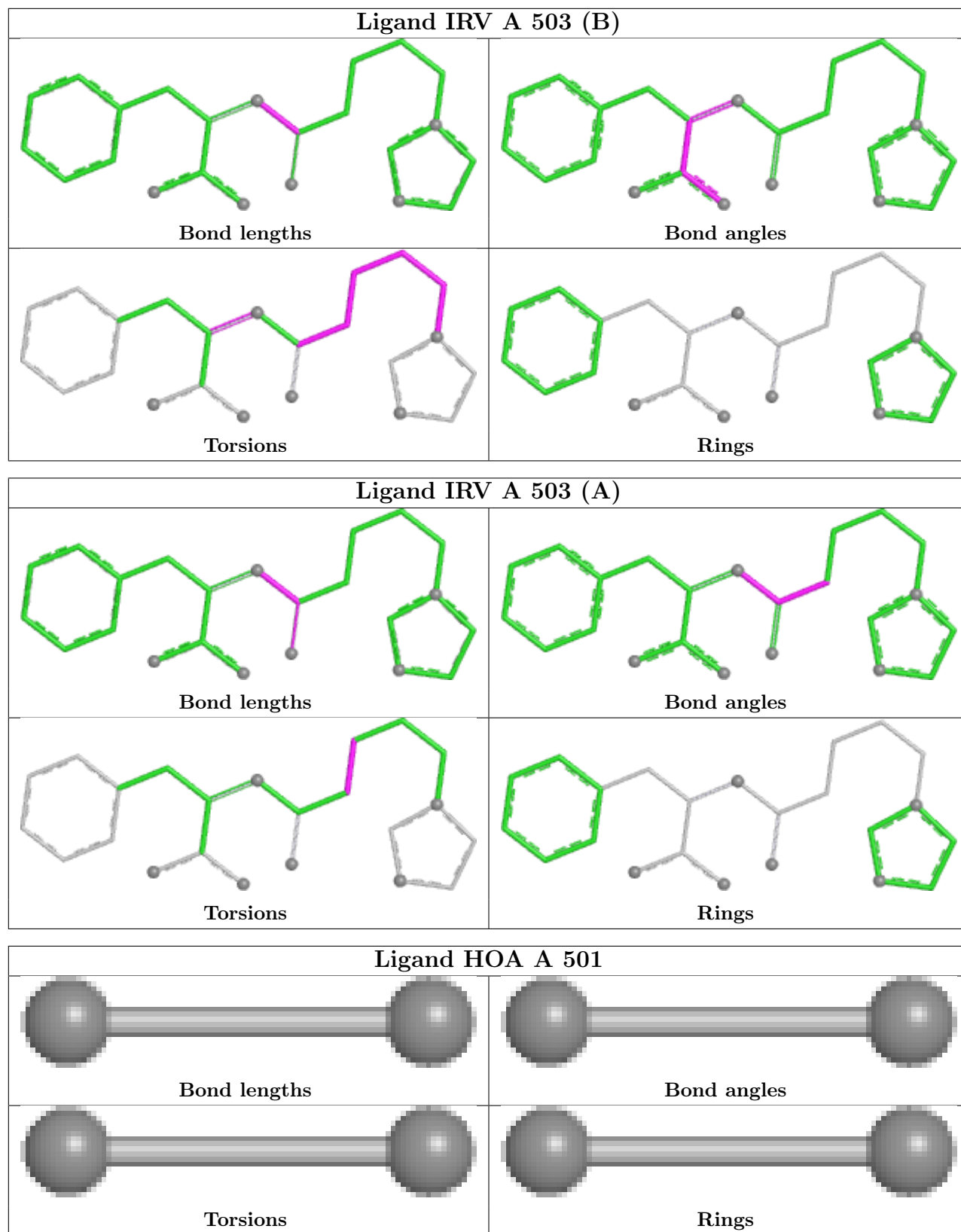
6 monomers are involved in 9 short contacts:

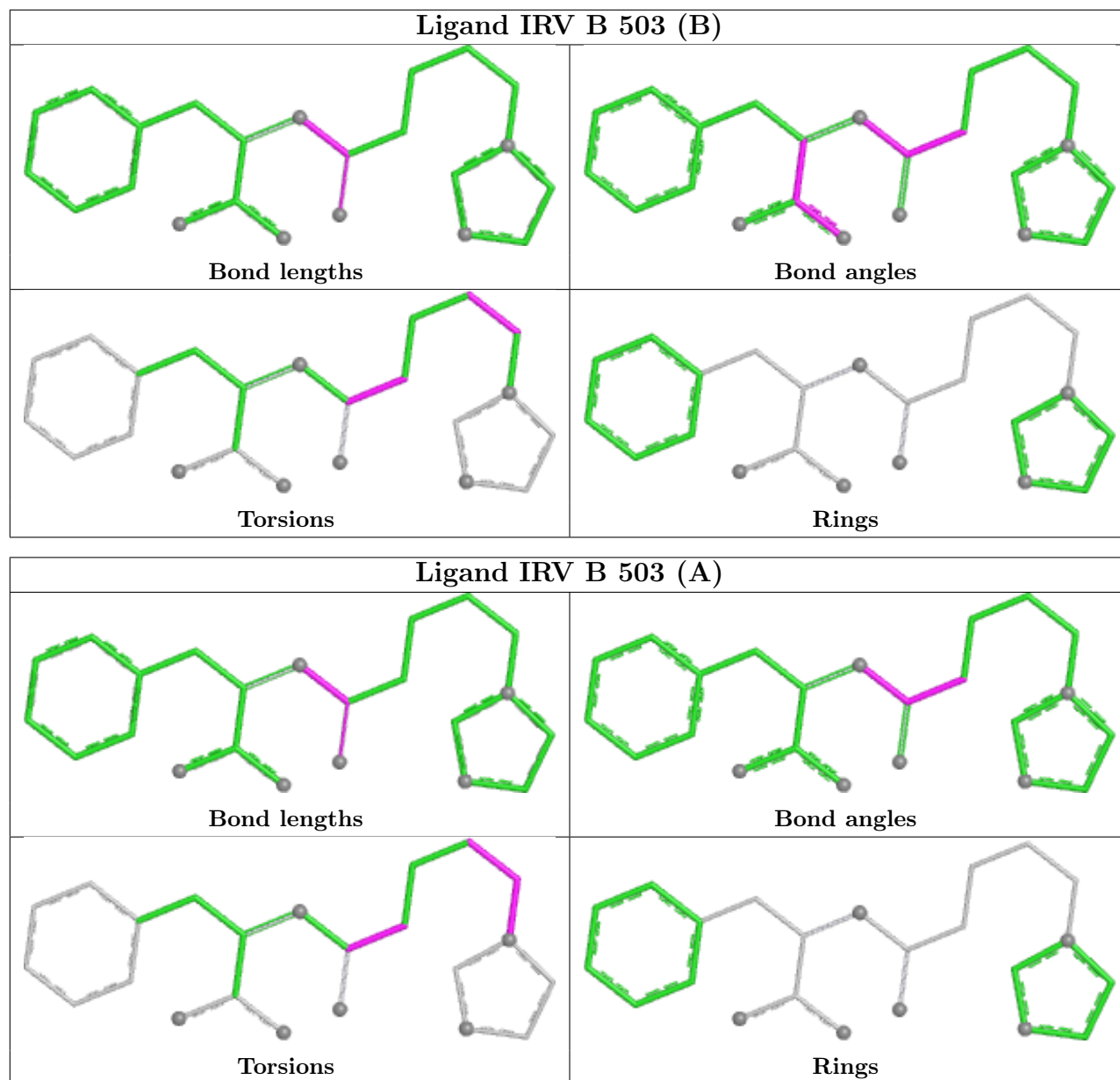
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	502	HOA	1	0
3	B	501	HEM	3	0
4	A	503[B]	IRV	1	0
4	A	503[A]	IRV	2	0
4	B	503[B]	IRV	1	0
3	A	502	HEM	2	0

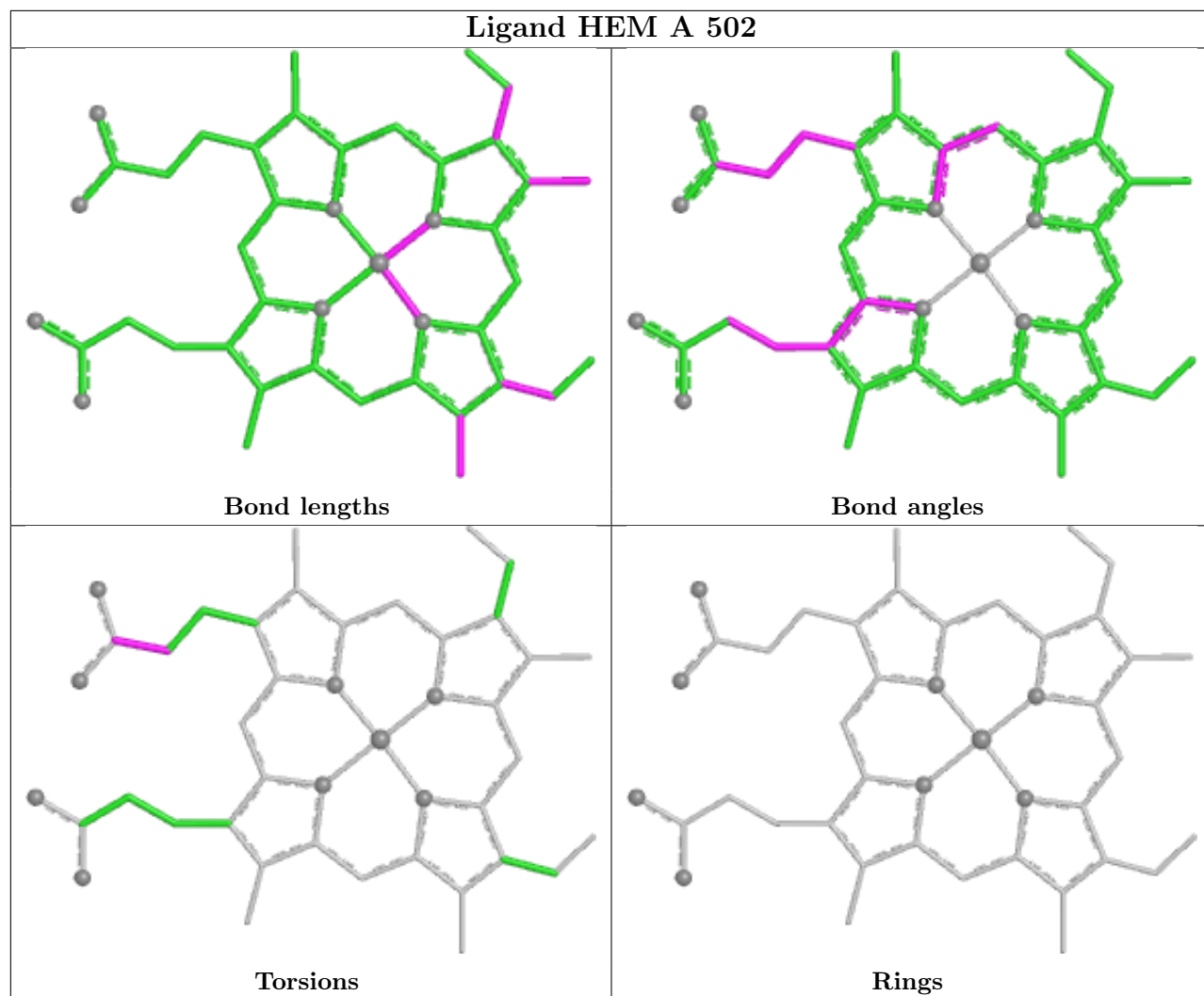
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

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	451/466 (96%)	0.22	16 (3%) 47 47	12, 33, 70, 99	1 (0%)
1	B	451/466 (96%)	0.26	23 (5%) 33 32	12, 34, 74, 112	1 (0%)
All	All	902/932 (96%)	0.24	39 (4%) 40 39	12, 33, 70, 112	2 (0%)

All (39) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	382	PRO	4.5
1	B	383	SER	4.0
1	B	381	ASN	3.7
1	A	197	ALA	3.7
1	B	215	LEU	3.7
1	A	437	LEU	3.6
1	A	371	VAL	3.5
1	A	136	ASP	3.4
1	A	196	PRO	3.3
1	A	189	GLN	3.1
1	B	194	ASP	3.1
1	A	231	ASP	3.1
1	B	191	ALA	3.0
1	B	384	ALA	2.9
1	B	193	PRO	2.9
1	A	191	ALA	2.9
1	B	192	ASN	2.8
1	B	189	GLN	2.8
1	A	199	ASP	2.8
1	A	195	ASP	2.7
1	B	168	ASP	2.7
1	B	136	ASP	2.7
1	B	197	ALA	2.7
1	A	302	VAL	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	198	TYR	2.6
1	B	369	ASP	2.5
1	B	385	ILE	2.5
1	A	193	PRO	2.5
1	B	196	PRO	2.5
1	B	169	GLN	2.4
1	A	225	ALA	2.4
1	B	198	TYR	2.3
1	B	219	ILE	2.3
1	B	45	PRO	2.1
1	A	211	VAL	2.1
1	B	366	ILE	2.1
1	B	218	LYS	2.0
1	A	158	PHE	2.0
1	B	199	ASP	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](#)

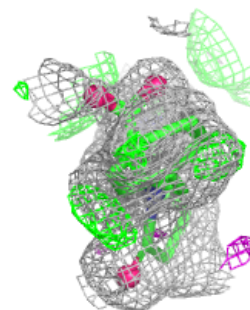
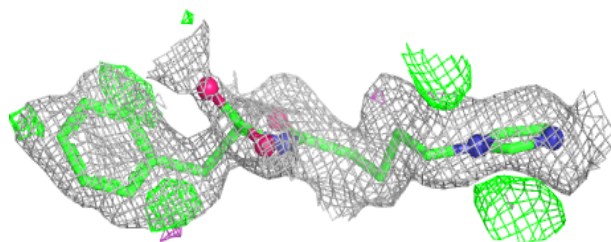
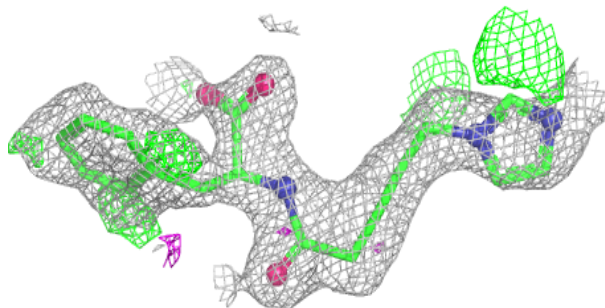
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	IRV	B	503[A]	23/23	0.77	0.17	26,41,65,66	23
4	IRV	B	503[B]	23/23	0.77	0.17	25,53,65,66	23
4	IRV	A	503[A]	23/23	0.82	0.16	22,43,54,55	23
4	IRV	A	503[B]	23/23	0.82	0.16	20,36,55,56	23
2	HOA	B	502	2/2	0.89	0.17	31,31,31,48	0
2	HOA	A	501	2/2	0.93	0.12	34,34,34,44	0
3	HEM	A	502	43/43	0.99	0.05	19,25,30,48	0
3	HEM	B	501	43/43	0.99	0.05	17,24,28,39	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.

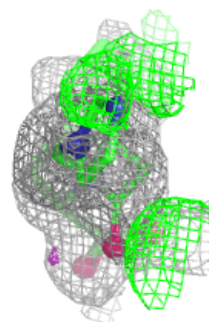
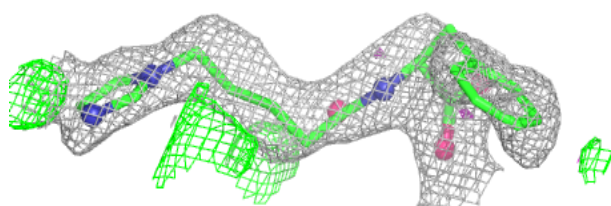
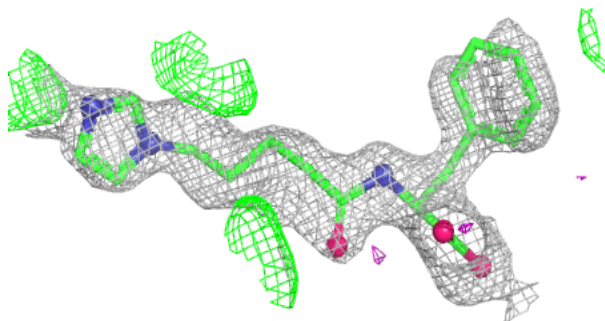
Electron density around IRV B 503 (A):

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

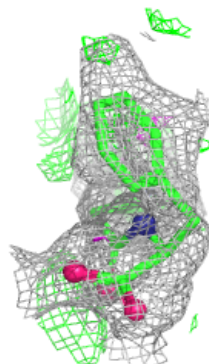
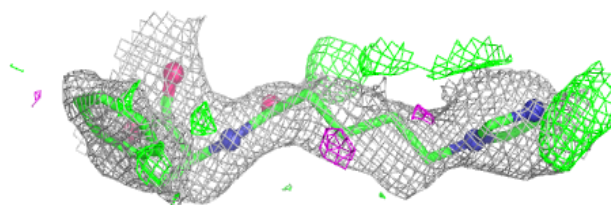
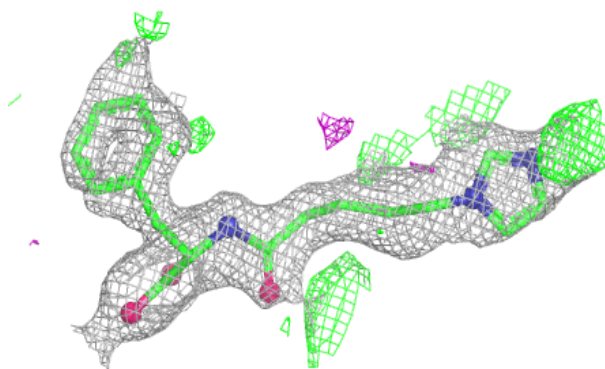


Electron density around IRV B 503 (B):

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

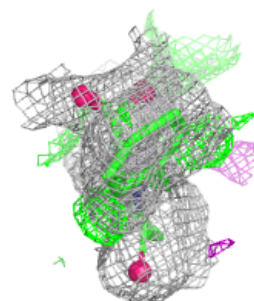
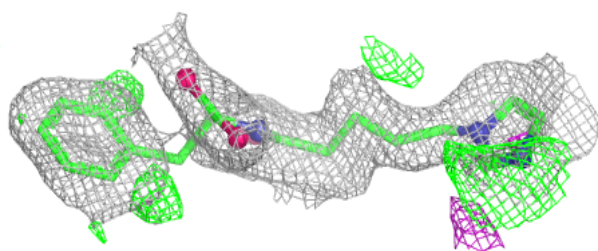
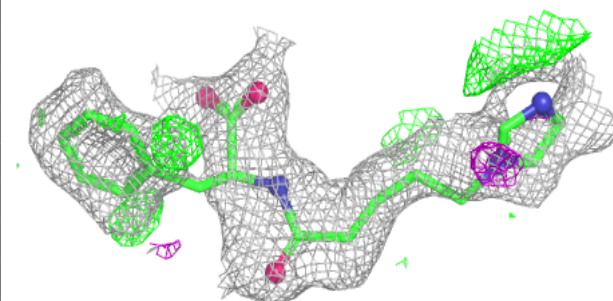
**Electron density around IRV A 503 (A):**

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

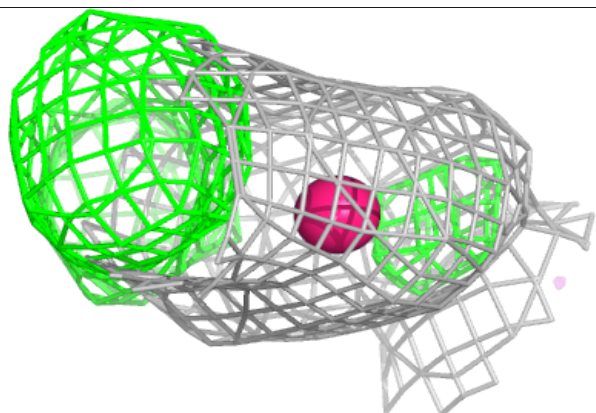
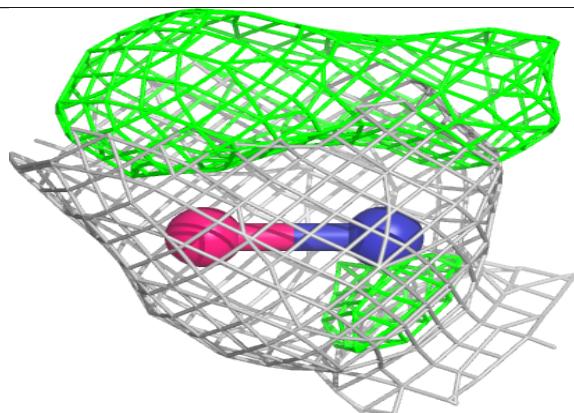
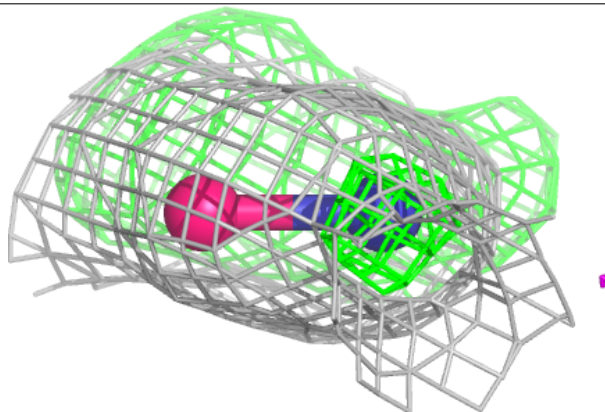


Electron density around IRV A 503 (B):

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and green (positive)

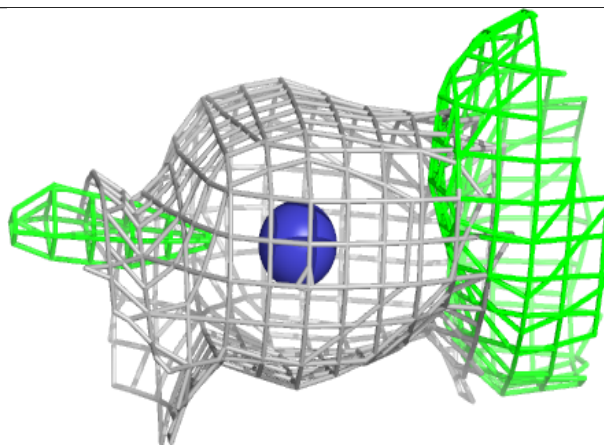
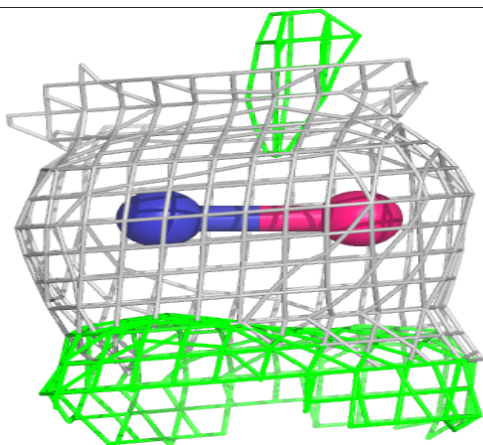
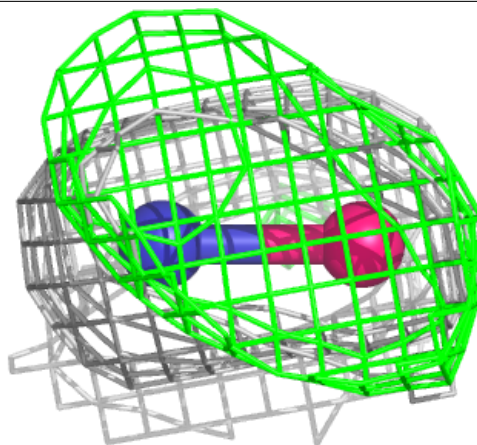
**Electron density around HOA B 502:**

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



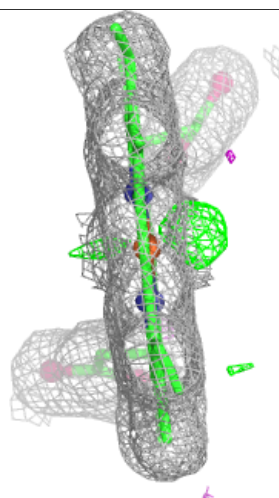
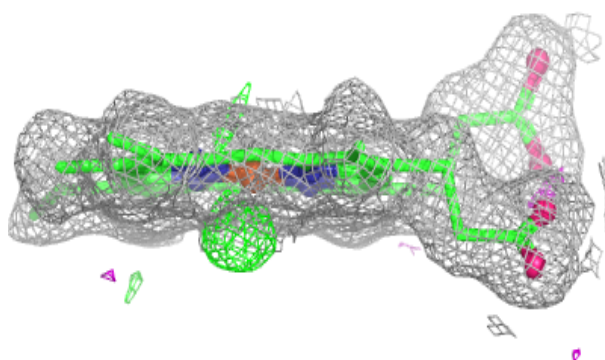
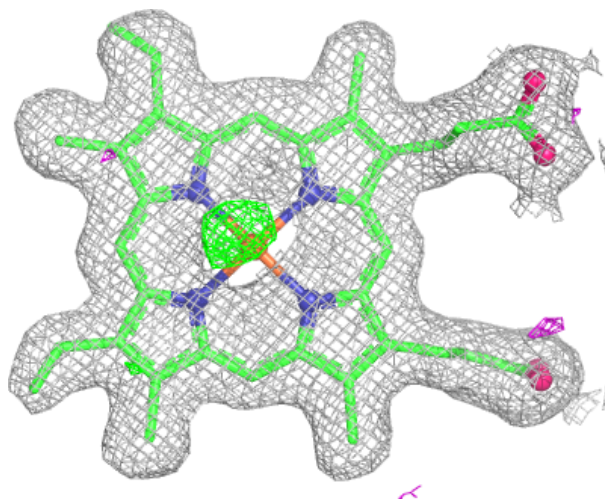
Electron density around HOA A 501:

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and green (positive)



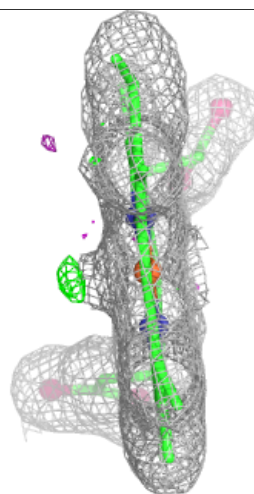
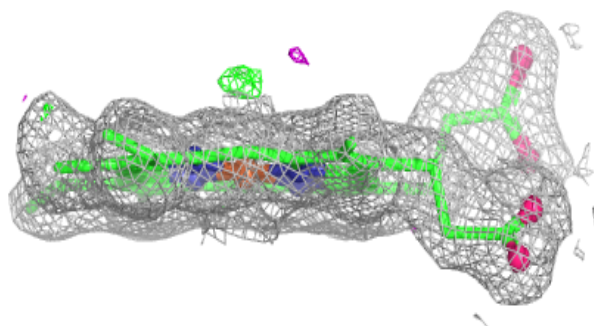
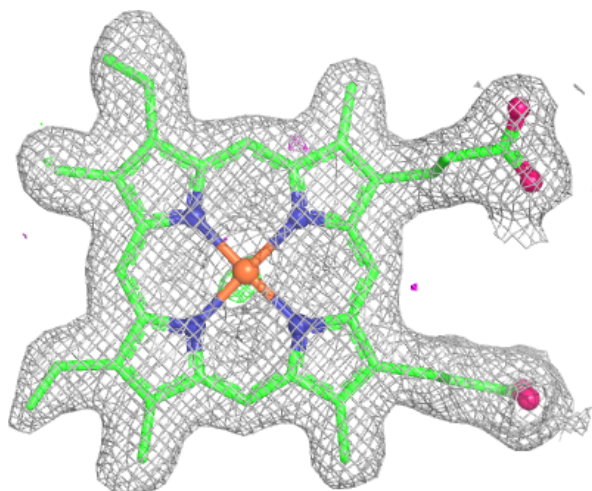
Electron density around HEM A 502:

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



Electron density around HEM 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)



6.5 Other polymers ⓘ

There are no such residues in this entry.