



Full wwPDB X-ray Structure Validation Report ⓘ

Jul 16, 2026 – 08:10 AM JST

PDB ID : 9L7U / pdb_0000917u
Title : Crystal structure of P450 BM3 F87A/V78S/L75N mutant complex with (-)-Ambroxide
Authors : Dong, S.; Feng, Y.G.
Deposited on : 2024-12-27
Resolution : 2.70 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	:	1.8.5 (274361), CSD as541be (2020)
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	:	9.0.015 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.50

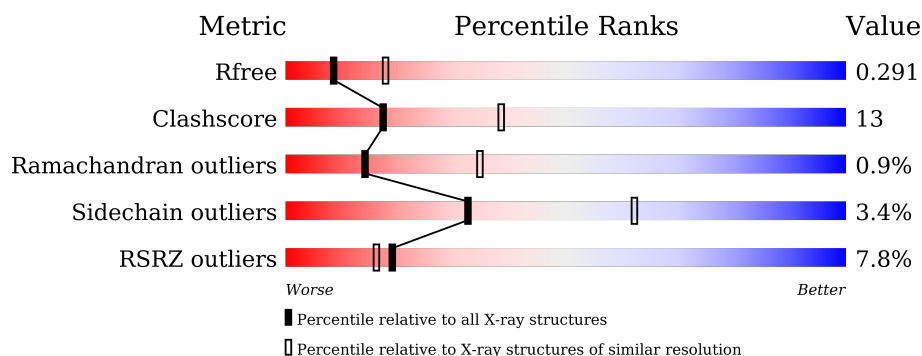
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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	456	8% 72% 25% .
1	B	456	7% 75% 22% .

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.

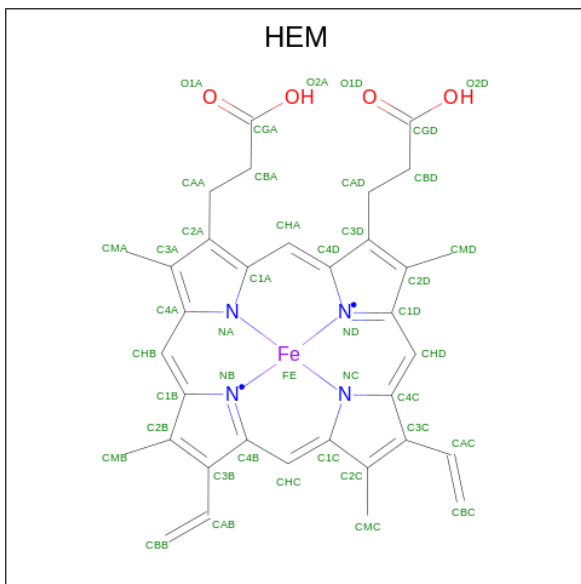
- Molecule 1 is a protein called Bifunctional cytochrome P450/NADPH-P450 reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	455	Total 3616	C 2306	N 618	O 675	S 17	0	0	0
1	B	455	Total 3622	C 2311	N 618	O 676	S 17	0	0	0

There are 6 discrepancies between the modelled and reference sequences:

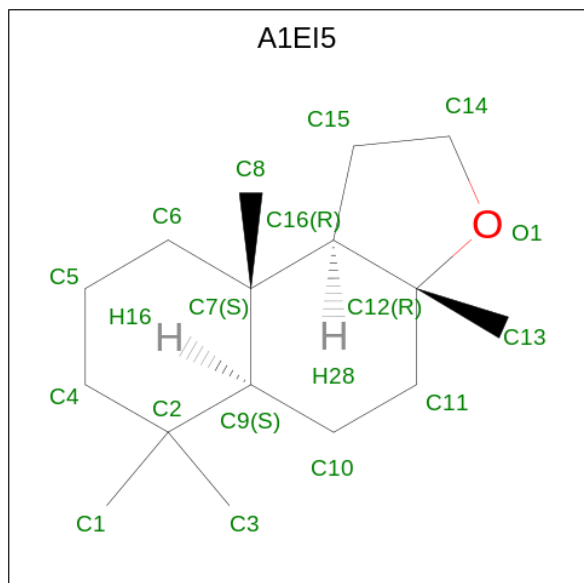
Chain	Residue	Modelled	Actual	Comment	Reference
A	75	ASN	LEU	engineered mutation	UNP F2Q7T0
A	78	SER	VAL	engineered mutation	UNP F2Q7T0
A	87	ALA	PHE	engineered mutation	UNP F2Q7T0
B	75	ASN	LEU	engineered mutation	UNP F2Q7T0
B	78	SER	VAL	engineered mutation	UNP F2Q7T0
B	87	ALA	PHE	engineered mutation	UNP F2Q7T0

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



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

- Molecule 3 is (3 {a} {R},5 {a} {S},9 {a} {S},9 {b} {R})-3 {a},6,6,9 {a}-tetramethyl-2,4,5,5 {a},7,8,9,9 {b}-octahydro-1 {H}-benzo[e][1]benzofuran (CCD ID: A1EI5) (formula: C₁₆H₂₈O) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O		
			17	16	1		
3	B	1	Total	C	O		
			17	16	1		

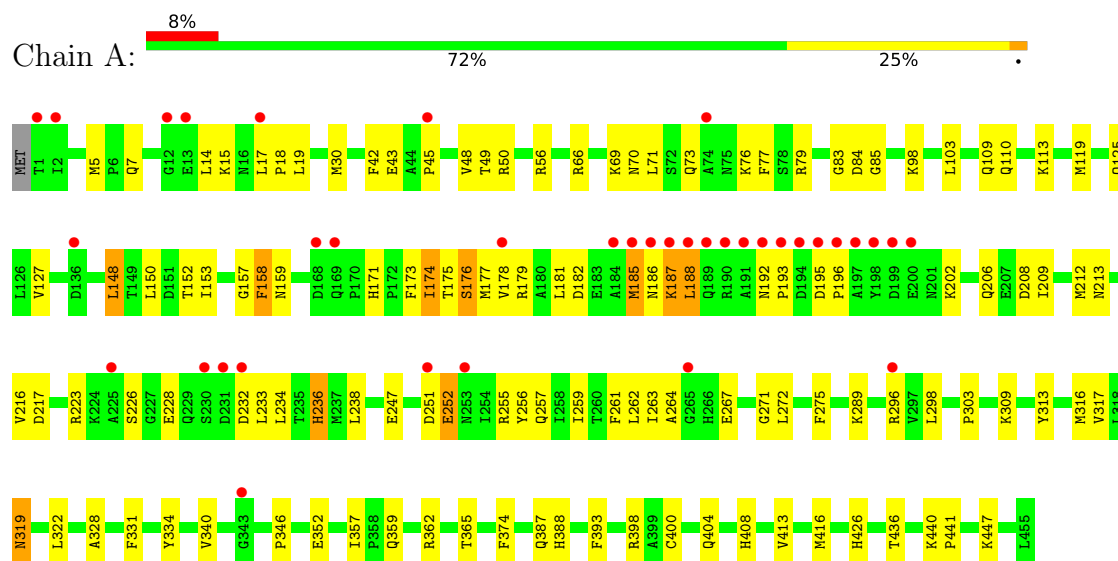
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	140	Total	O		
			140	140		
4	B	146	Total	O		
			146	146		

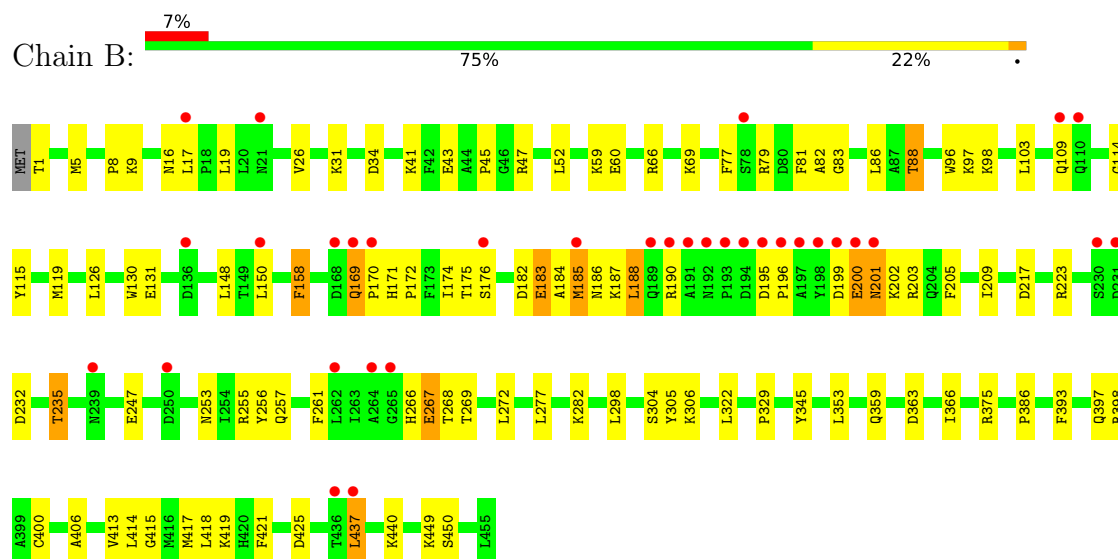
3 Residue-property plots

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: Bifunctional cytochrome P450/NADPH-P450 reductase



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4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	57.17Å 158.22Å 58.72Å 90.00° 93.63° 90.00°	Depositor
Resolution (Å)	79.11 – 2.70 79.11 – 2.70	Depositor EDS
% Data completeness (in resolution range)	98.8 (79.11-2.70) 90.7 (79.11-2.70)	Depositor EDS
R_{merge}	0.33	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.06 (at 2.69Å)	Xtriage
Refinement program	PHENIX (1.20.1_4487: ???)	Depositor
R, R_{free}	0.231 , 0.275 0.251 , 0.291	Depositor DCC
R_{free} test set	2044 reflections (7.21%)	wwPDB-VP
Wilson B-factor (Å ²)	19.7	Xtriage
Anisotropy	0.341	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 43.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.42$, $\langle L^2 \rangle = 0.24$	Xtriage
Estimated twinning fraction	0.056 for l,-k,h	Xtriage
F_o, F_c correlation	0.84	EDS
Total number of atoms	7644	wwPDB-VP
Average B, all atoms (Å ²)	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.25% 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: A1EI5, 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	0.21	0/3697	0.39	0/4999
1	B	0.28	0/3703	0.39	0/5006
All	All	0.25	0/7400	0.39	0/10005

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	3616	0	3559	95	0
1	B	3622	0	3562	82	0
2	A	43	0	30	2	0
2	B	43	0	30	3	0
3	A	17	0	0	0	0
3	B	17	0	0	2	0
4	A	140	0	0	17	0
4	B	146	0	0	16	0
All	All	7644	0	7181	182	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 13.

All (182) 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:200:GLU:HG2	1:B:203:ARG:NH2	1.85	0.92
1:A:177:MET:CE	1:A:262:LEU:HB3	2.05	0.86
1:A:71:LEU:O	1:A:76:LYS:HE3	1.76	0.85
1:A:171:HIS:O	1:A:175:THR:HG23	1.75	0.85
1:B:414:LEU:O	4:B:601:HOH:O	1.96	0.82
1:B:150:LEU:HD13	1:B:174:ILE:HD13	1.60	0.82
1:A:109:GLN:HG2	1:A:404:GLN:HG2	1.62	0.82
1:A:182:ASP:O	1:A:186:ASN:HB2	1.78	0.81
1:A:148:LEU:O	1:A:152:THR:OG1	2.00	0.79
1:A:252:GLU:OE2	4:A:601:HOH:O	2.00	0.79
1:B:148:LEU:HD11	1:B:413:VAL:HG21	1.64	0.77
1:B:200:GLU:OE1	4:B:602:HOH:O	2.03	0.75
1:A:83:GLY:HA3	1:A:256:TYR:CZ	2.23	0.74
1:B:202:LYS:HA	1:B:205:PHE:HB3	1.69	0.73
1:A:171:HIS:HB3	1:A:174:ILE:HD12	1.69	0.73
1:A:216:VAL:CG1	1:A:255:ARG:HG3	2.19	0.73
1:A:150:LEU:HD13	1:A:174:ILE:HG12	1.70	0.72
1:B:306:LYS:NZ	4:B:605:HOH:O	2.21	0.72
1:A:159:ASN:ND2	4:A:606:HOH:O	2.21	0.72
1:A:228:GLU:OE2	4:A:602:HOH:O	2.07	0.72
1:A:185:MET:O	1:A:185:MET:HE2	1.90	0.70
1:A:177:MET:HE1	1:A:262:LEU:HB3	1.71	0.70
1:B:98:LYS:HE2	1:B:247:GLU:HB2	1.75	0.69
1:B:185:MET:HE2	1:B:437:LEU:HD13	1.75	0.69
1:A:83:GLY:HA3	1:A:256:TYR:CE2	2.29	0.68
1:B:17:LEU:HD22	1:B:45:PRO:HD2	1.74	0.68
1:A:272:LEU:HD13	1:A:322:LEU:HG	1.77	0.67
1:A:17:LEU:HD22	1:A:45:PRO:HD2	1.78	0.65
1:B:5:MET:HE2	1:B:41:LYS:HB2	1.79	0.65
1:A:264:ALA:O	4:A:603:HOH:O	2.15	0.65
1:A:171:HIS:HB3	1:A:174:ILE:CD1	2.27	0.64
1:A:436:THR:N	4:A:615:HOH:O	2.27	0.64
1:A:216:VAL:HG12	1:A:255:ARG:HG3	1.79	0.64
1:A:264:ALA:HB1	4:A:603:HOH:O	1.97	0.64
1:A:98:LYS:HE3	1:A:247:GLU:HB2	1.79	0.63
1:A:362:ARG:NH1	4:A:612:HOH:O	2.30	0.63
1:A:66:ARG:NH1	4:A:622:HOH:O	2.32	0.62
1:B:47:ARG:NH2	4:B:623:HOH:O	2.31	0.62
1:A:296:ARG:HG2	1:A:296:ARG:O	1.98	0.62
1:A:217:ASP:OD1	1:A:255:ARG:HD3	2.01	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:183:GLU:O	1:B:187:LYS:HG2	2.00	0.61
1:B:34:ASP:OD1	1:B:359:GLN:NE2	2.34	0.60
1:B:375:ARG:NH2	4:B:625:HOH:O	2.34	0.60
2:B:501:HEM:HBC2	2:B:501:HEM:HMC2	1.83	0.60
1:B:171:HIS:CG	1:B:172:PRO:HD2	2.37	0.60
1:B:66:ARG:NH1	4:B:620:HOH:O	2.31	0.59
1:A:179:ARG:HG2	1:A:208:ASP:OD2	2.01	0.59
1:A:209:ILE:O	1:A:213:ASN:ND2	2.25	0.59
1:B:277:LEU:HD13	1:B:417:MET:HE2	1.85	0.59
1:B:59:LYS:NZ	4:B:628:HOH:O	2.36	0.58
2:A:501:HEM:HMC2	2:A:501:HEM:HBC2	1.84	0.58
1:B:200:GLU:O	1:B:203:ARG:N	2.29	0.58
1:A:125:GLN:NE2	4:A:626:HOH:O	2.33	0.57
2:A:501:HEM:HBB2	2:A:501:HEM:HMB2	1.87	0.57
1:A:177:MET:HE2	1:A:262:LEU:HB3	1.86	0.57
1:A:56:ARG:NH1	4:A:625:HOH:O	2.33	0.56
1:A:340:VAL:HA	1:A:346:PRO:HA	1.87	0.56
1:B:272:LEU:HD13	1:B:322:LEU:HG	1.87	0.56
1:B:205:PHE:O	1:B:209:ILE:HG12	2.06	0.56
1:A:153:ILE:O	1:A:157:GLY:N	2.32	0.55
1:B:9:LYS:HE3	1:B:19:LEU:HD11	1.89	0.55
1:A:176:SER:OG	1:A:208:ASP:OD1	2.23	0.55
1:B:182:ASP:O	1:B:186:ASN:HB2	2.05	0.55
1:B:200:GLU:C	1:B:202:LYS:N	2.64	0.55
1:B:366:ILE:N	4:B:631:HOH:O	2.36	0.55
1:A:157:GLY:HA2	1:A:233:LEU:HB2	1.88	0.54
1:B:282:LYS:NZ	1:B:425:ASP:OD2	2.36	0.54
1:B:366:ILE:HD12	1:B:386:PRO:HG3	1.90	0.54
1:B:69:LYS:HB3	1:B:398:ARG:HG3	1.90	0.53
1:A:42:PHE:HB3	1:A:49:THR:HG22	1.91	0.53
1:B:232:ASP:OD1	1:B:235:THR:OG1	2.20	0.53
1:A:251:ASP:HB3	4:A:729:HOH:O	2.07	0.53
1:B:269:THR:HG23	1:B:406:ALA:HB1	1.91	0.52
1:A:110:GLN:HG2	1:A:113:LYS:HE3	1.91	0.52
1:A:79:ARG:NH1	4:A:641:HOH:O	2.42	0.52
1:B:397:GLN:NE2	4:B:640:HOH:O	2.40	0.52
1:B:200:GLU:HG2	1:B:203:ARG:HH22	1.68	0.52
3:B:502:A1EI5:C4	4:B:604:HOH:O	2.58	0.52
1:B:60:GLU:HG2	1:B:66:ARG:HH12	1.74	0.51
1:B:9:LYS:O	1:B:16:ASN:ND2	2.43	0.51
1:B:223:ARG:NH2	1:B:232:ASP:OD2	2.44	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:232:ASP:OD2	1:A:234:LEU:HB3	2.11	0.51
1:A:267:GLU:OE2	4:A:604:HOH:O	2.19	0.51
1:A:5:MET:HE1	1:A:50:ARG:HH11	1.76	0.50
1:A:173:PHE:HE1	1:A:177:MET:HE3	1.76	0.50
1:B:217:ASP:OD2	1:B:255:ARG:NH1	2.44	0.50
1:B:304:SER:N	4:B:608:HOH:O	2.39	0.50
1:B:77:PHE:CE2	1:B:187:LYS:O	2.65	0.50
1:B:115:TYR:HB2	1:B:119:MET:HE3	1.93	0.50
1:A:5:MET:HE1	1:A:50:ARG:HD3	1.92	0.49
1:A:298:LEU:HD22	1:A:303:PRO:HB3	1.93	0.49
1:B:130:TRP:CH2	1:B:417:MET:HE3	2.47	0.49
1:B:393:PHE:HB3	1:B:400:CYS:HB3	1.92	0.49
1:A:158:PHE:HA	1:A:232:ASP:OD2	2.12	0.49
1:A:69:LYS:HB3	1:A:398:ARG:HG3	1.94	0.49
1:B:130:TRP:HH2	1:B:417:MET:HE3	1.78	0.48
1:A:73:GLN:HA	1:A:73:GLN:OE1	2.13	0.48
1:B:8:PRO:HD2	1:B:41:LYS:O	2.13	0.48
1:B:200:GLU:C	1:B:202:LYS:H	2.21	0.48
1:B:415:GLY:C	4:B:601:HOH:O	2.57	0.48
2:B:501:HEM:HMB1	2:B:501:HEM:HBB2	1.95	0.48
1:A:202:LYS:O	1:A:206:GLN:HG2	2.13	0.48
1:B:267:GLU:OE2	1:B:440:LYS:NZ	2.45	0.48
1:A:127:VAL:HG21	1:A:416:MET:HE2	1.94	0.48
1:B:184:ALA:O	1:B:188:LEU:HD23	2.13	0.48
1:A:352:GLU:OE2	4:A:605:HOH:O	2.20	0.47
1:B:200:GLU:O	1:B:202:LYS:N	2.47	0.47
1:B:109:GLN:O	1:B:305:TYR:OH	2.33	0.47
1:B:81:PHE:HB3	1:B:209:ILE:HD12	1.97	0.47
1:B:437:LEU:HD23	3:B:502:A1EI5:C11	2.44	0.47
1:A:84:ASP:OD1	1:A:256:TYR:CD2	2.68	0.47
1:A:289:LYS:NZ	4:A:628:HOH:O	2.33	0.47
1:B:200:GLU:CG	1:B:203:ARG:NH2	2.68	0.47
1:A:223:ARG:HH12	1:A:232:ASP:CG	2.24	0.46
1:A:393:PHE:HB3	1:A:400:CYS:HB3	1.98	0.46
1:A:236:HIS:ND1	1:A:236:HIS:N	2.62	0.46
1:B:158:PHE:CD1	1:B:158:PHE:N	2.84	0.46
1:A:187:LYS:HB2	1:A:187:LYS:HE3	1.63	0.46
1:A:226:SER:OG	1:A:228:GLU:HG3	2.16	0.46
1:B:52:LEU:HD11	1:B:353:LEU:HD13	1.96	0.46
1:B:449:LYS:HD2	1:B:450:SER:N	2.31	0.46
1:A:85:GLY:HA2	1:A:257:GLN:NE2	2.31	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:289:LYS:HE3	1:A:313:TYR:CZ	2.51	0.45
1:A:158:PHE:CD1	1:A:158:PHE:N	2.85	0.45
1:A:174:ILE:O	1:A:178:VAL:HG13	2.17	0.45
1:A:309:LYS:HD3	1:A:309:LYS:HA	1.60	0.45
1:B:103:LEU:HD13	1:B:261:PHE:HZ	1.81	0.45
1:B:298:LEU:HB2	1:B:419:LYS:HD2	1.99	0.45
1:B:77:PHE:CD1	1:B:187:LYS:HE3	2.52	0.45
1:B:79:ARG:HD2	1:B:83:GLY:O	2.17	0.45
1:B:158:PHE:N	1:B:158:PHE:HD1	2.14	0.45
1:A:176:SER:OG	1:A:208:ASP:HA	2.17	0.45
1:A:216:VAL:HG11	1:A:255:ARG:HG3	1.95	0.44
1:B:31:LYS:HA	1:B:31:LYS:HD3	1.86	0.44
1:B:199:ASP:C	1:B:201:ASN:N	2.76	0.44
1:B:114:GLY:N	4:B:654:HOH:O	2.50	0.44
1:B:223:ARG:HH22	1:B:232:ASP:CG	2.26	0.44
1:A:319:ASN:HD22	1:A:319:ASN:HA	1.56	0.44
1:A:14:LEU:HB3	1:A:18:PRO:HD3	2.00	0.44
1:A:70:ASN:HB2	1:A:334:TYR:HB3	1.99	0.44
1:A:275:PHE:CD1	1:A:441:PRO:HG3	2.53	0.44
1:B:131:GLU:HG2	1:B:421:PHE:CZ	2.53	0.44
1:B:418:LEU:N	4:B:601:HOH:O	2.03	0.44
1:A:238:LEU:HD23	1:A:238:LEU:HA	1.86	0.43
1:B:26:VAL:HG11	1:B:329:PRO:HB2	2.00	0.43
1:B:268:THR:HB	2:B:501:HEM:C3B	2.53	0.43
1:A:103:LEU:HD13	1:A:261:PHE:HZ	1.83	0.43
1:B:363:ASP:HB3	4:B:714:HOH:O	2.19	0.43
1:A:119:MET:SD	1:A:408:HIS:HD2	2.42	0.43
1:A:30:MET:HG2	1:A:359:GLN:HG2	2.00	0.43
1:A:259:ILE:HG22	1:A:263:ILE:HD11	1.99	0.43
1:A:125:GLN:NE2	4:A:646:HOH:O	2.45	0.43
1:B:171:HIS:HB3	1:B:174:ILE:HG13	2.01	0.43
1:B:253:ASN:O	1:B:257:GLN:HG2	2.18	0.43
1:A:177:MET:O	1:A:181:LEU:HG	2.18	0.43
1:A:19:LEU:HD12	1:A:19:LEU:HA	1.86	0.42
1:B:5:MET:HB3	1:B:345:TYR:CE1	2.54	0.42
1:B:175:THR:OG1	4:B:603:HOH:O	2.09	0.42
1:A:252:GLU:CD	4:A:601:HOH:O	2.57	0.42
1:A:271:GLY:CA	1:A:440:LYS:HG2	2.49	0.42
1:B:83:GLY:HA3	1:B:256:TYR:CE2	2.54	0.42
1:A:426:HIS:CD2	1:A:447:LYS:HE2	2.55	0.42
1:A:328:ALA:O	1:A:357:ILE:HD12	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:188:LEU:HD22	1:B:188:LEU:HA	1.89	0.42
1:B:199:ASP:C	1:B:201:ASN:H	2.26	0.42
1:B:169:GLN:HG2	1:B:170:PRO:HD2	2.00	0.42
1:A:73:GLN:NE2	1:A:77:PHE:CZ	2.88	0.42
1:A:109:GLN:OE1	1:A:309:LYS:HE3	2.20	0.42
1:A:313:TYR:HA	1:A:316:MET:HB2	2.01	0.42
1:B:86:LEU:HD23	1:B:86:LEU:HA	1.92	0.42
1:A:317:VAL:HG13	1:A:374:PHE:HZ	1.85	0.41
1:B:150:LEU:HG	1:B:266:HIS:HB2	2.02	0.41
1:A:148:LEU:HD11	1:A:413:VAL:HG21	2.01	0.41
1:A:43:GLU:HG2	1:A:48:VAL:HG22	2.03	0.41
1:A:15:LYS:HA	1:A:15:LYS:HD2	1.80	0.41
1:A:173:PHE:CZ	1:A:212:MET:HG2	2.56	0.41
1:A:387:GLN:HG2	1:A:388:HIS:CD2	2.56	0.41
1:B:96:TRP:CZ3	1:B:97:LYS:HD3	2.56	0.41
1:A:185:MET:HB2	1:A:185:MET:HE3	1.65	0.41
1:A:188:LEU:HD22	1:A:188:LEU:HA	1.84	0.41
1:B:82:ALA:HB3	1:B:88:THR:HG21	2.03	0.40
1:A:331:PHE:HD1	1:A:357:ILE:HD11	1.87	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	453/456 (99%)	436 (96%)	13 (3%)	4 (1%)	14	35
1	B	453/456 (99%)	439 (97%)	10 (2%)	4 (1%)	14	35
All	All	906/912 (99%)	875 (97%)	23 (2%)	8 (1%)	14	35

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	192	ASN
1	A	193	PRO
1	A	195	ASP
1	A	196	PRO
1	B	195	ASP
1	B	196	PRO
1	B	190	ARG
1	B	201	ASN

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	385/399 (96%)	373 (97%)	12 (3%)	35	65
1	B	385/399 (96%)	371 (96%)	14 (4%)	31	60
All	All	770/798 (96%)	744 (97%)	26 (3%)	32	62

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	GLN
1	A	148	LEU
1	A	158	PHE
1	A	174	ILE
1	A	176	SER
1	A	185	MET
1	A	187	LYS
1	A	188	LEU
1	A	236	HIS
1	A	252	GLU
1	A	319	ASN
1	A	365	THR
1	B	1	THR
1	B	43	GLU
1	B	88	THR
1	B	126	LEU

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Mol	Chain	Res	Type
1	B	158	PHE
1	B	169	GLN
1	B	176	SER
1	B	183	GLU
1	B	185	MET
1	B	188	LEU
1	B	200	GLU
1	B	235	THR
1	B	267	GLU
1	B	437	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	201	ASN
1	A	319	ASN
1	B	204	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](#)

4 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the

expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	A1EI5	A	502	-	19,19,19	0.21	0	30,32,32	1.04	1 (3%)
2	HEM	B	501	4	50,50,50	1.44	8 (16%)	66,82,82	1.23	8 (12%)
2	HEM	A	501	4	50,50,50	1.43	8 (16%)	66,82,82	1.24	8 (12%)
3	A1EI5	B	502	-	19,19,19	0.21	0	30,32,32	1.01	1 (3%)

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	A1EI5	A	502	-	-	-	0/3/3/3
2	HEM	B	501	4	-	0/14/54/54	-
2	HEM	A	501	4	-	0/14/54/54	-
3	A1EI5	B	502	-	-	-	0/3/3/3

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	HEM	FE-NB	3.63	2.06	1.94
2	B	501	HEM	FE-NB	3.46	2.05	1.94
2	A	501	HEM	FE-NC	3.42	2.06	1.95
2	B	501	HEM	FE-NC	3.40	2.06	1.95
2	B	501	HEM	FE-ND	3.30	2.05	1.94
2	B	501	HEM	FE-NA	3.28	2.06	1.95
2	A	501	HEM	FE-ND	3.10	2.04	1.94
2	B	501	HEM	CAB-C3B	3.08	1.55	1.47
2	A	501	HEM	CAB-C3B	3.07	1.55	1.47
2	A	501	HEM	CAC-C3C	3.06	1.55	1.47
2	B	501	HEM	CAC-C3C	3.02	1.55	1.47
2	A	501	HEM	FE-NA	2.91	2.05	1.95
2	A	501	HEM	CMC-C2C	2.05	1.55	1.50
2	B	501	HEM	CMC-C2C	2.04	1.55	1.50
2	B	501	HEM	CMB-C2B	2.03	1.55	1.50
2	A	501	HEM	CMB-C2B	2.02	1.55	1.50

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	502	A1EI5	O1-C12-C16	-3.92	95.66	101.78
3	B	502	A1EI5	O1-C12-C16	-3.68	96.03	101.78
2	A	501	HEM	C4A-NA-C1A	3.15	108.43	105.35
2	A	501	HEM	C4D-ND-C1D	3.14	108.32	105.07
2	B	501	HEM	C4A-NA-C1A	2.98	108.27	105.35
2	B	501	HEM	C1B-NB-C4B	2.97	108.14	105.07
2	A	501	HEM	C1B-NB-C4B	2.95	108.12	105.07
2	B	501	HEM	C4D-ND-C1D	2.95	108.12	105.07
2	B	501	HEM	C4C-NC-C1C	2.63	107.93	105.35
2	A	501	HEM	C4C-NC-C1C	2.56	107.85	105.35
2	A	501	HEM	C2A-C1A-NA	-2.32	107.55	110.15
2	B	501	HEM	C3B-C2B-C1B	2.31	108.20	106.49
2	A	501	HEM	C3D-C4D-ND	-2.21	107.71	110.17
2	B	501	HEM	CAD-CBD-CGD	-2.19	108.90	113.60
2	A	501	HEM	C3B-C2B-C1B	2.18	108.11	106.49
2	B	501	HEM	C2A-C1A-NA	-2.11	107.78	110.15
2	A	501	HEM	CAD-CBD-CGD	-2.11	109.06	113.60
2	B	501	HEM	C3D-C4D-ND	-2.10	107.82	110.17

There are no chirality outliers.

There are no torsion outliers.

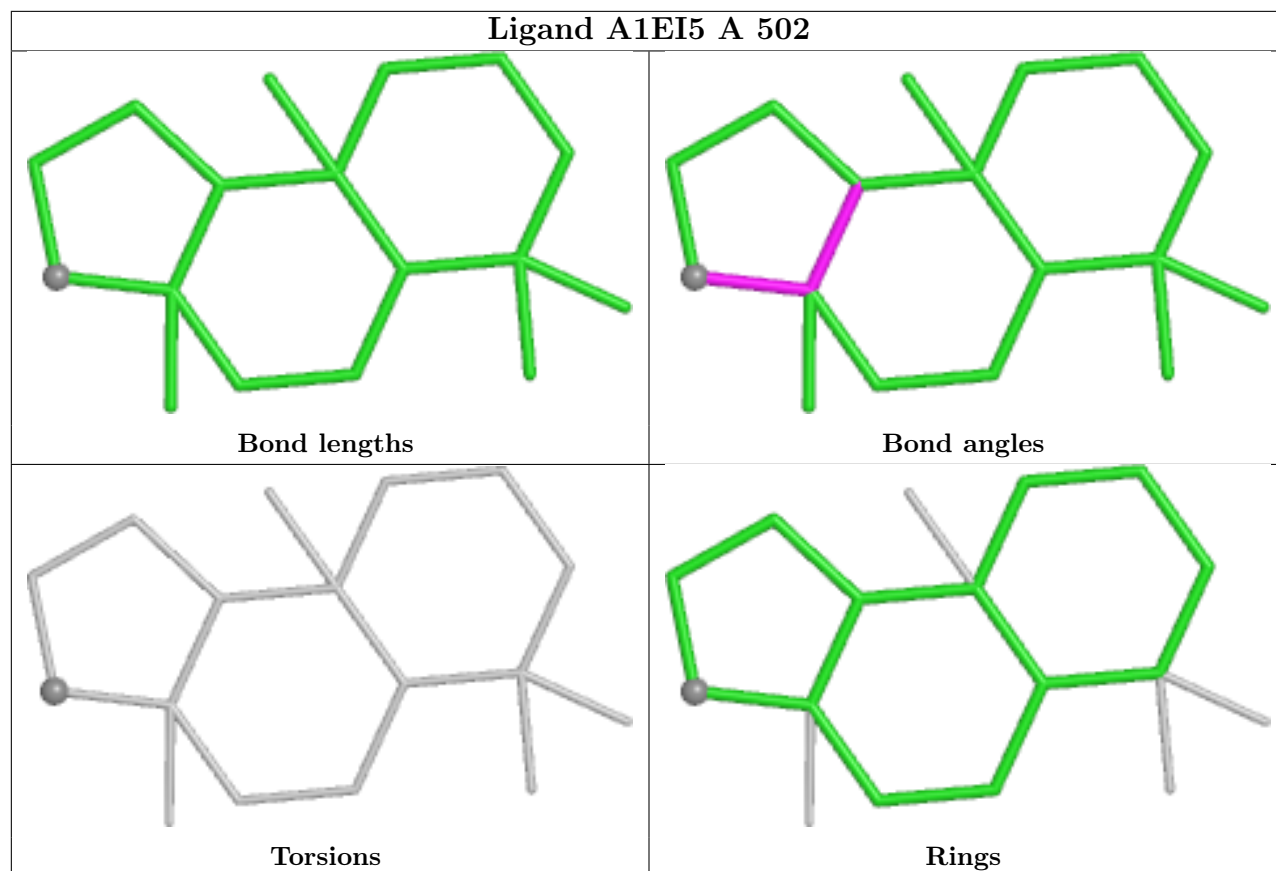
There are no ring outliers.

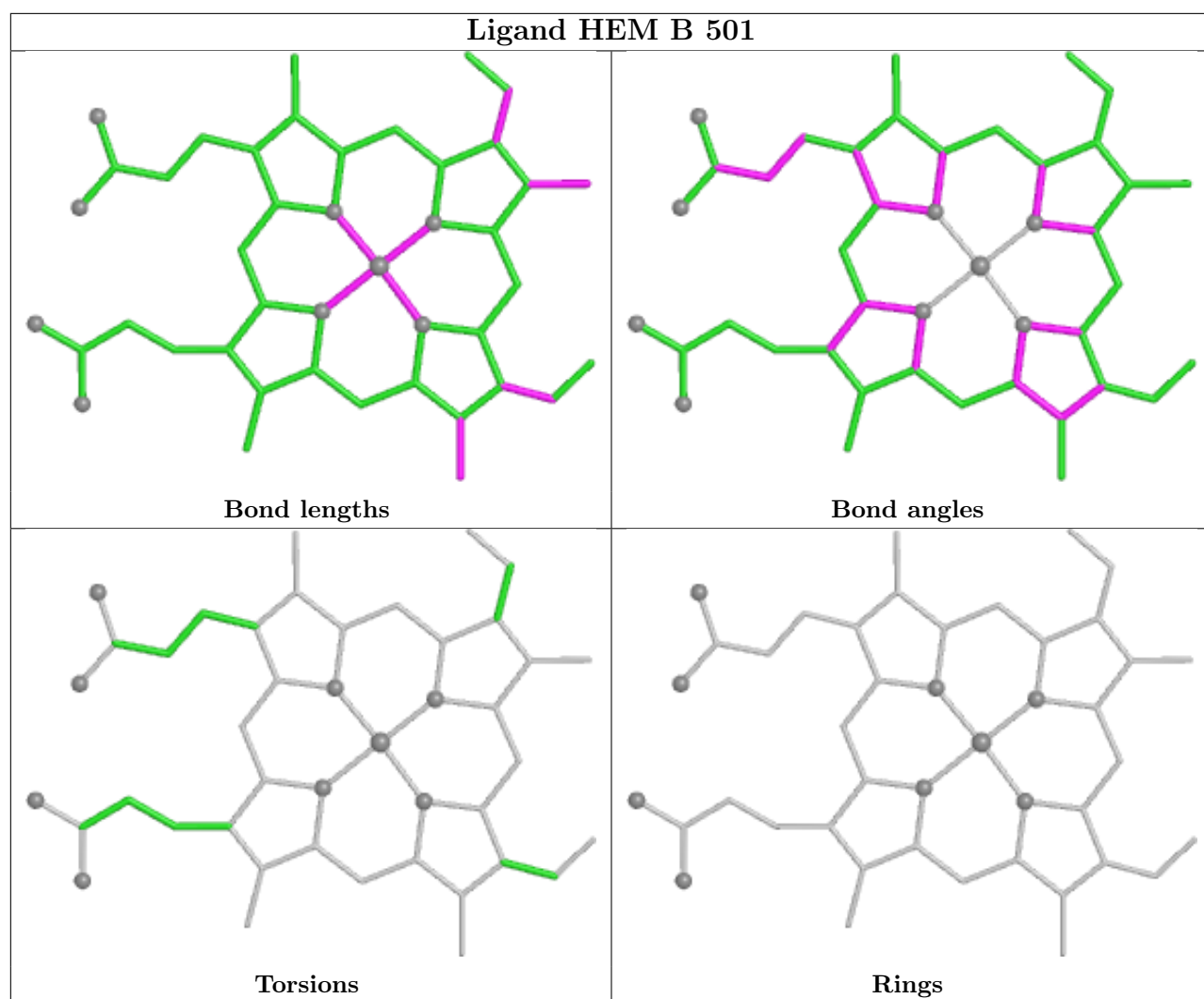
3 monomers are involved in 7 short contacts:

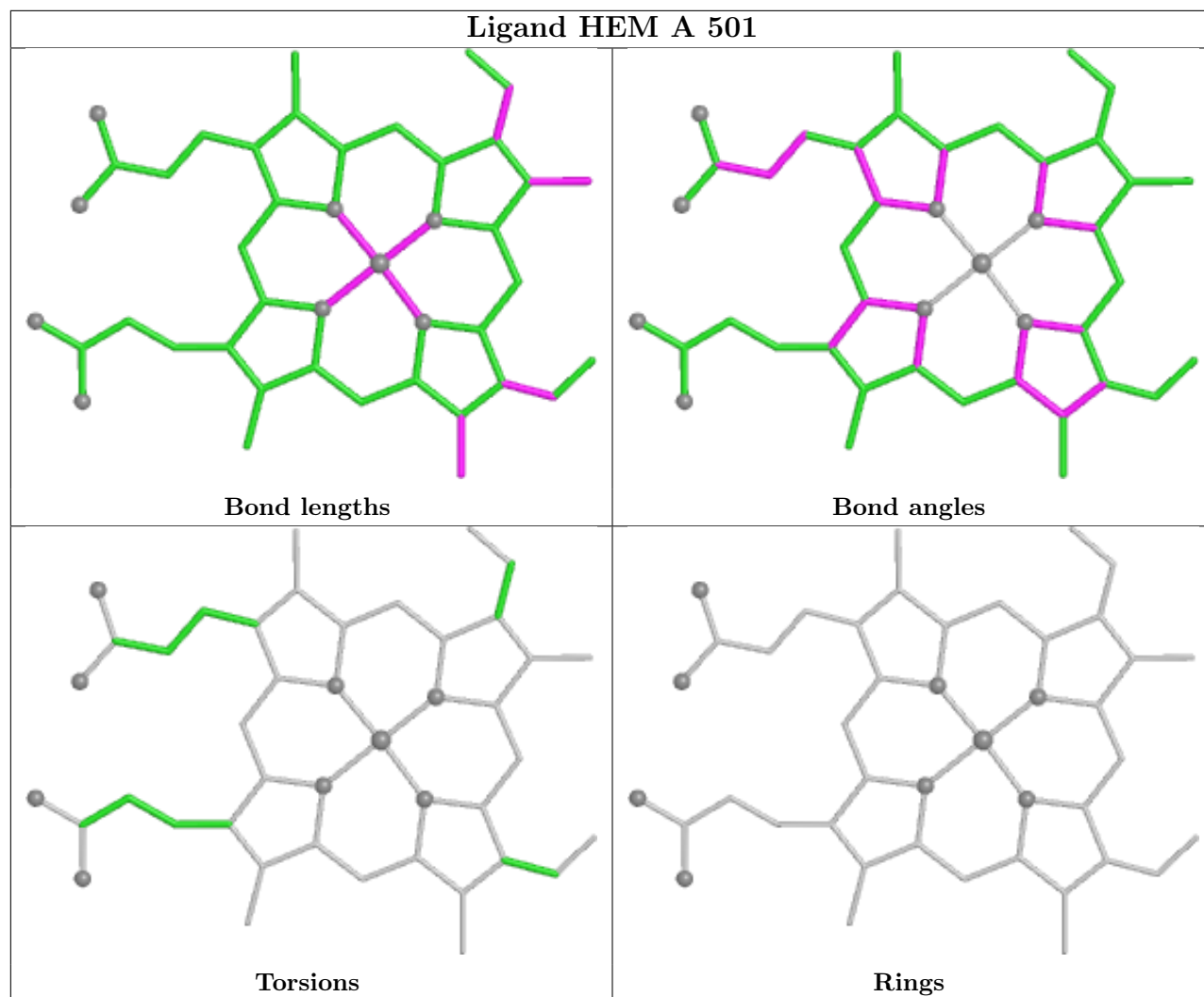
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	501	HEM	3	0
2	A	501	HEM	2	0
3	B	502	A1EI5	2	0

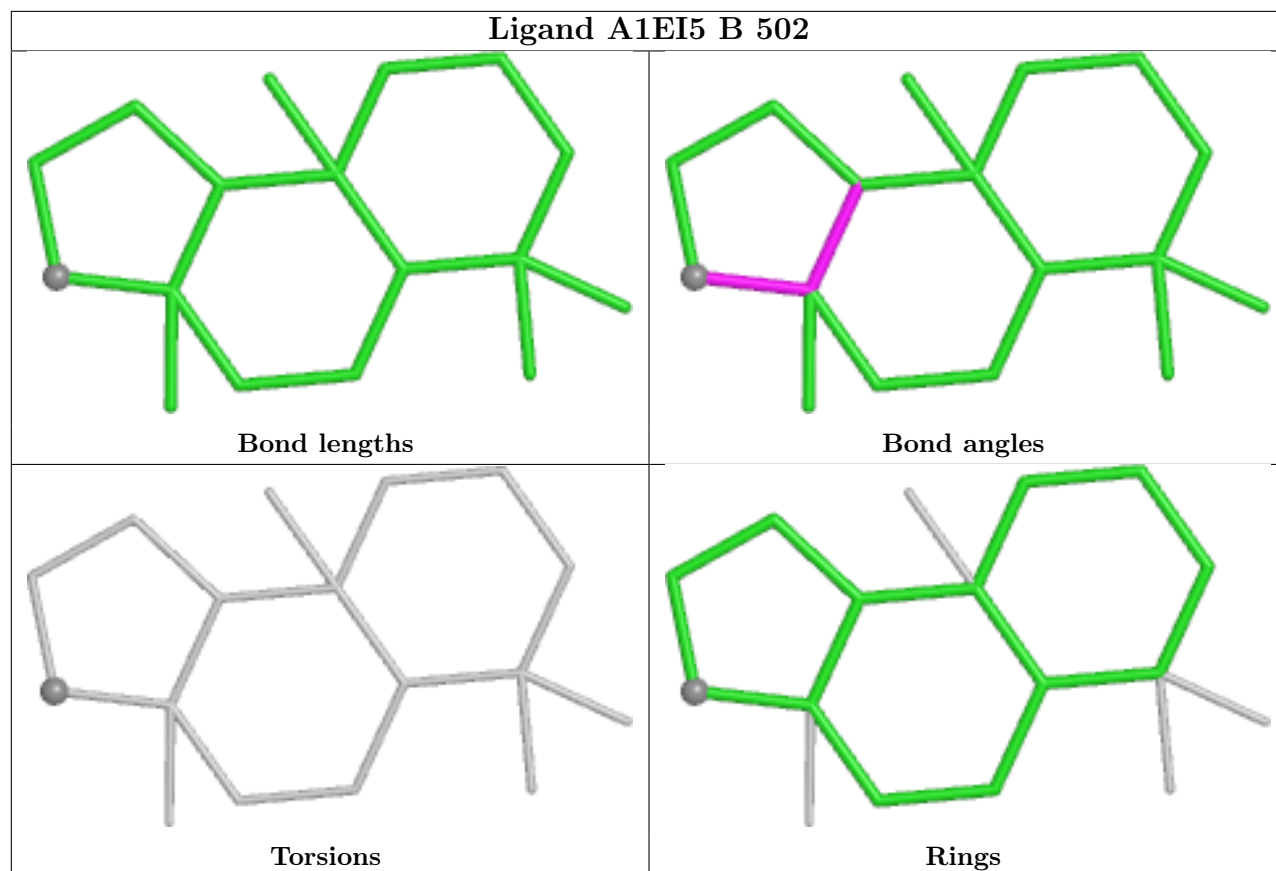
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

Ligand A1EI5 A 502









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	455/456 (99%)	0.85	37 (8%) 18 15	13, 24, 49, 74	0
1	B	455/456 (99%)	0.82	34 (7%) 20 17	15, 23, 43, 66	0
All	All	910/912 (99%)	0.83	71 (7%) 19 16	13, 23, 47, 74	0

All (71) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	191	ALA	12.5
1	B	196	PRO	8.4
1	B	192	ASN	6.2
1	A	198	TYR	6.0
1	B	198	TYR	5.9
1	A	191	ALA	5.8
1	B	189	GLN	5.7
1	B	199	ASP	5.6
1	A	194	ASP	5.2
1	A	190	ARG	5.2
1	B	194	ASP	5.1
1	A	195	ASP	5.0
1	B	195	ASP	4.9
1	A	192	ASN	4.9
1	A	197	ALA	4.8
1	A	188	LEU	4.8
1	A	193	PRO	4.6
1	A	185	MET	4.5
1	A	253	ASN	4.4
1	A	187	LYS	4.4
1	A	186	ASN	4.1
1	A	196	PRO	4.1
1	B	136	ASP	4.0
1	A	230	SER	3.9

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Mol	Chain	Res	Type	RSRZ
1	A	169	GLN	3.9
1	A	168	ASP	3.8
1	A	1	THR	3.8
1	B	190	ARG	3.4
1	A	231	ASP	3.3
1	A	189	GLN	3.3
1	A	17	LEU	3.3
1	B	437	LEU	3.2
1	A	12	GLY	3.2
1	B	197	ALA	3.1
1	A	296	ARG	3.0
1	B	168	ASP	2.9
1	A	199	ASP	2.9
1	A	232	ASP	2.8
1	B	231	ASP	2.8
1	A	225	ALA	2.8
1	B	230	SER	2.7
1	A	200	GLU	2.7
1	B	193	PRO	2.7
1	B	200	GLU	2.6
1	A	136	ASP	2.6
1	A	184	ALA	2.6
1	B	109	GLN	2.6
1	A	45	PRO	2.5
1	A	343	GLY	2.5
1	B	17	LEU	2.5
1	B	110	GLN	2.4
1	A	2	ILE	2.4
1	B	239	ASN	2.4
1	B	21	ASN	2.4
1	A	251	ASP	2.4
1	B	185	MET	2.4
1	B	262	LEU	2.3
1	A	13	GLU	2.3
1	A	74	ALA	2.3
1	B	201	ASN	2.3
1	B	169	GLN	2.3
1	B	436	THR	2.3
1	A	265	GLY	2.3
1	B	264	ALA	2.2
1	B	250	ASP	2.2
1	A	178	VAL	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	176	SER	2.2
1	B	78	SER	2.1
1	B	265	GLY	2.1
1	B	170	PRO	2.0
1	B	150	LEU	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](#)

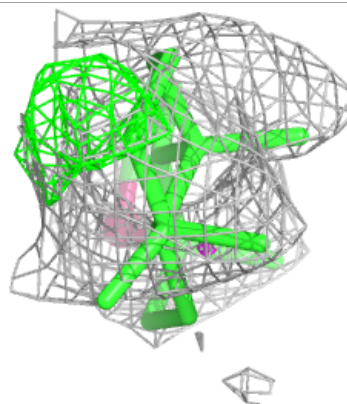
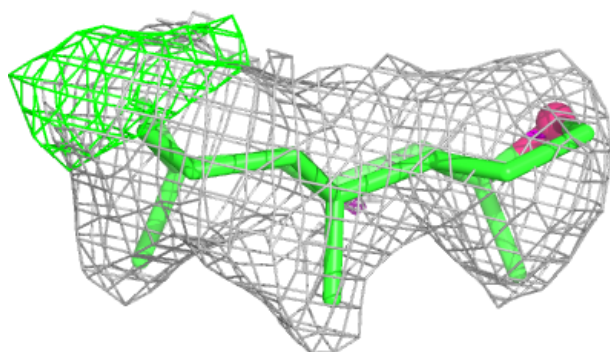
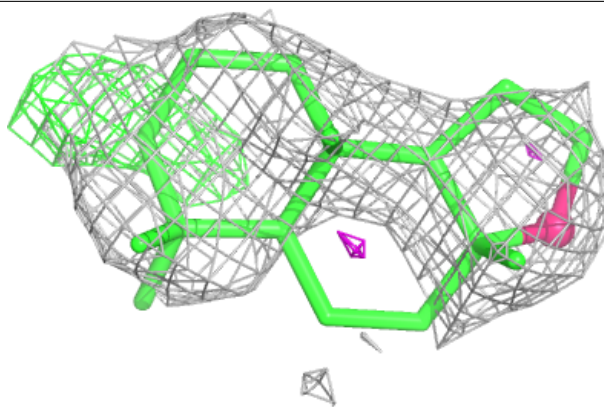
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	A1EI5	A	502	17/17	0.69	0.20	20,27,32,34	0
3	A1EI5	B	502	17/17	0.72	0.22	23,29,38,42	0
2	HEM	B	501	43/43	0.92	0.13	12,19,22,23	0
2	HEM	A	501	43/43	0.93	0.12	10,18,20,22	0

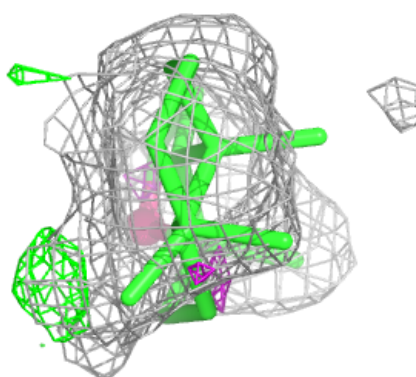
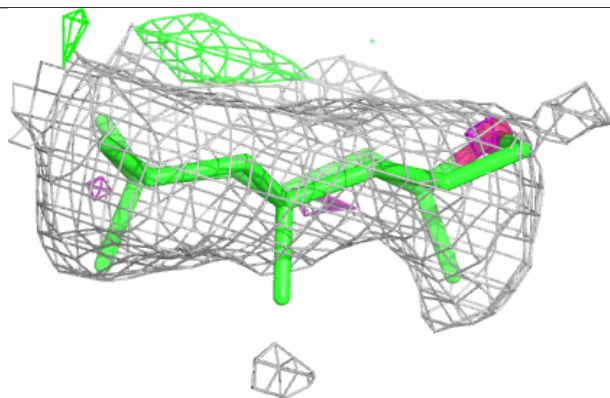
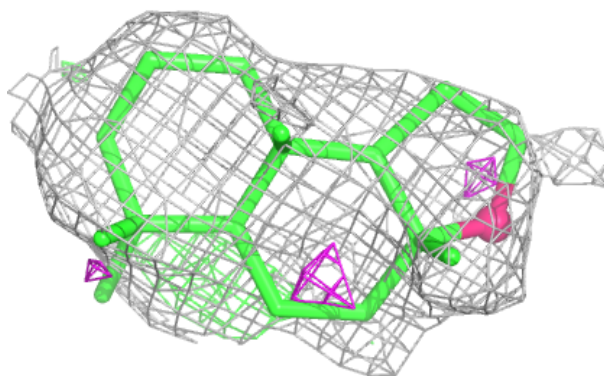
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around A1EI5 A 502:

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

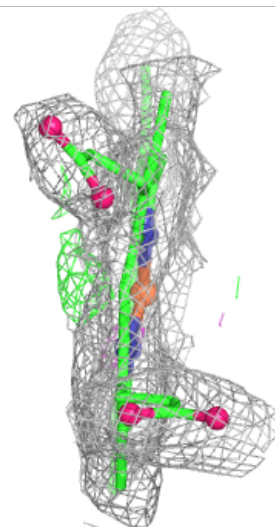
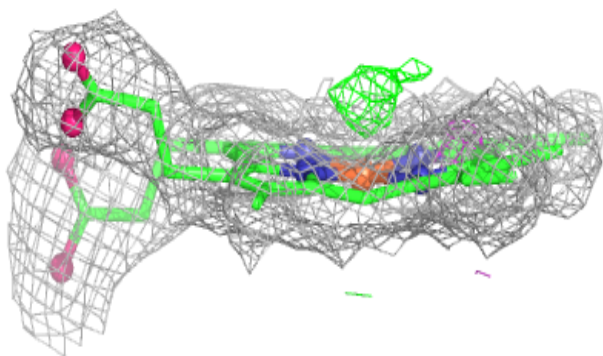
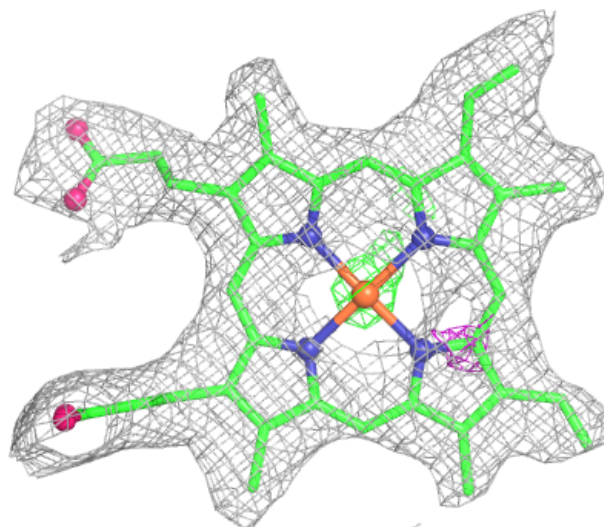
**Electron density around A1EI5 B 502:**

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



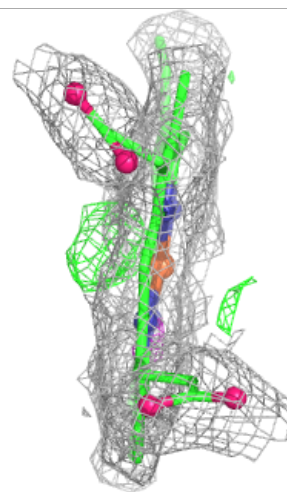
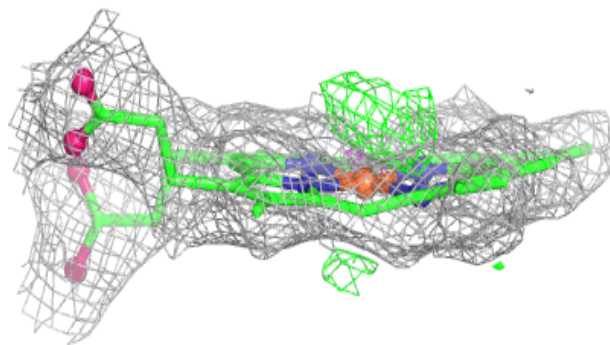
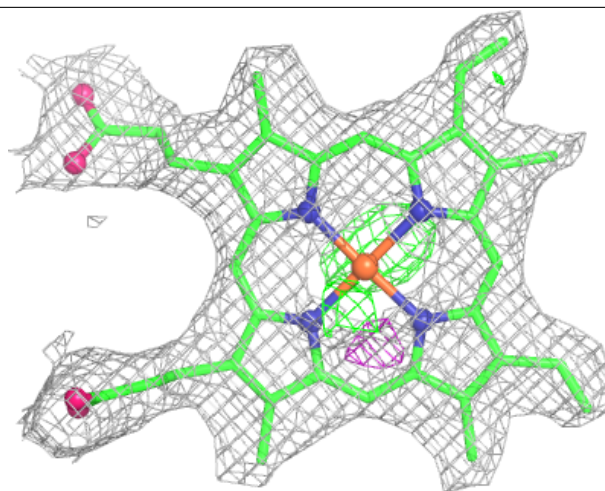
Electron density around HEM B 501:

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



Electron density around HEM A 501:

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



6.5 Other polymers [i](#)

There are no such residues in this entry.