



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 3, 2026 – 09:19 PM EDT

PDB ID : 1K2O / pdb_00001k2o
Title : Cytochrome P450Cam with Bound BIS(2,2'-BIPYRIDINE)-(5-METHYL-2-2'-BIPYRIDINE)-C2-ADAMANTANE RUTHENIUM (II)
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Deposited on : 2001-09-28
Resolution : 1.65 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)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.50

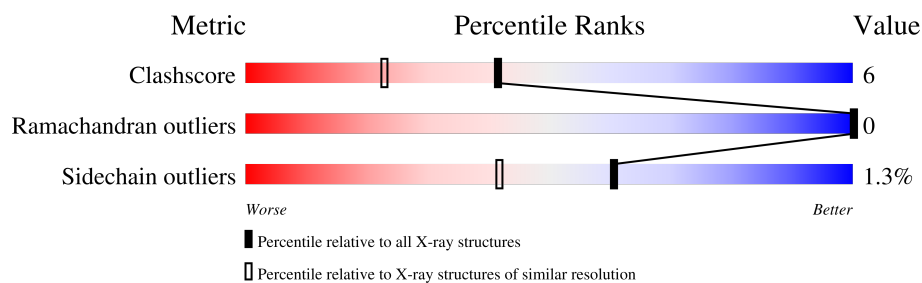
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.65 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))
Clashscore	190562	2662 (1.66-1.66)
Ramachandran outliers	187476	2621 (1.66-1.66)
Sidechain outliers	187428	2621 (1.66-1.66)

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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	414	
1	B	414	

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 7649 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.

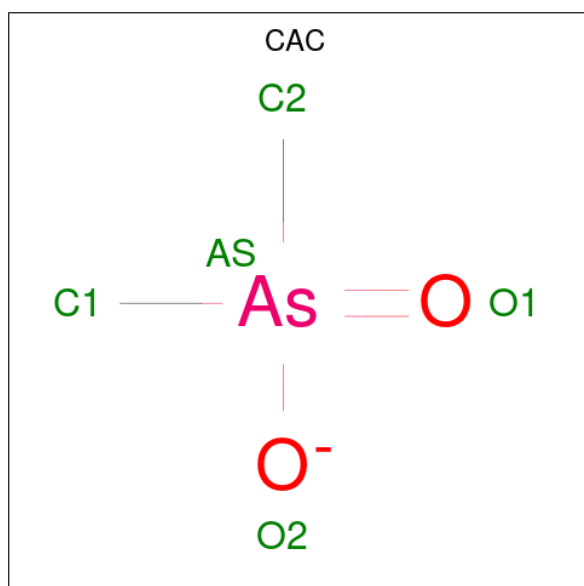
- Molecule 1 is a protein called Cytochrome P450CAM.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	406	Total	C	N	O	S	0	12	0
			3311	2094	579	620	18			
1	B	406	Total	C	N	O	S	0	6	0
			3258	2063	568	608	19			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	334	ALA	CYS	engineered mutation	UNP P00183
B	334	ALA	CYS	engineered mutation	UNP P00183

- Molecule 2 is CACODYLATE ION (CCD ID: CAC) (formula: $C_2H_6AsO_2$).



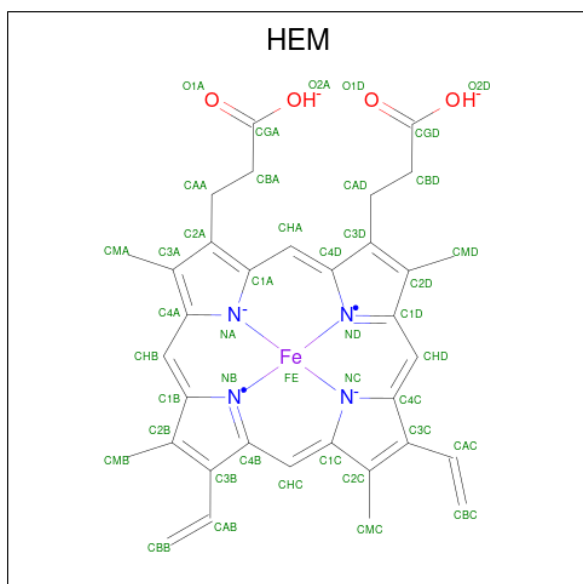
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	As	C	O	0	0
			4	1	2	1		

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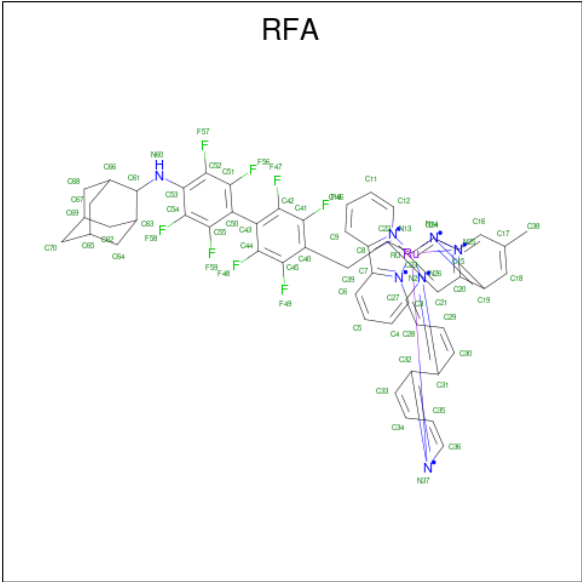
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	As	C	O	0	0
			4	1	2	1		
2	B	1	Total	As	C	O	0	0
			4	1	2	1		
2	B	1	Total	As	C	O	0	0
			4	1	2	1		
2	B	1	Total	As	C	O	0	0
			4	1	2	1		

- Molecule 3 is PROTOPORPHYRIN IX CONTAINING FE (CCD ID: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).



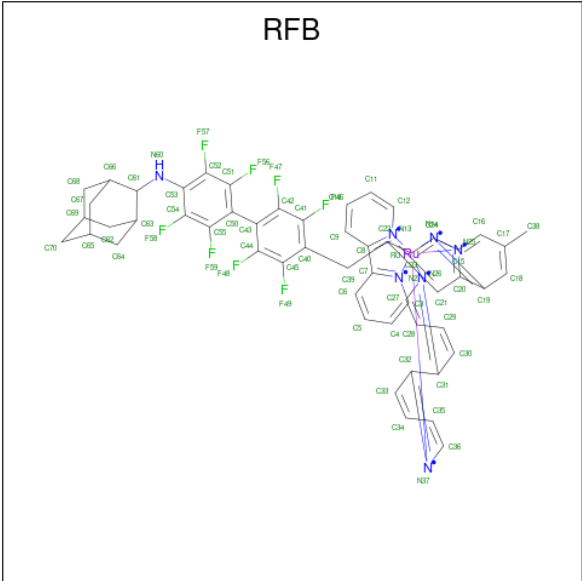
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	Fe	N	O	0
			43	34	1	4	4	
3	B	1	Total	C	Fe	N	O	0
			43	34	1	4	4	

- Molecule 4 is DELTA-BIS(2,2'-BIPYRIDINE)-(5-METHYL-2-2'-BIPYRIDINE)-C2-ADAMANTANE RUTHENIUM (II) (CCD ID: RFA) (formula: $C_{54}H_{43}F_8N_7Ru$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	F	N	Ru	0	1
			70	54	8	7	1		
4	B	1	Total	C	F	N	Ru	0	1
			70	54	8	7	1		

- Molecule 5 is LAMBDA-BIS(2,2'-BIPYRIDINE)-(5-METHYL-2-2'-BIPYRIDINE)-C2-AD AMANTANE RUTHENIUM (II) (CCD ID: RFB) (formula: $C_{54}H_{43}F_8N_7Ru$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total	C	F	N	Ru	0	1
			70	54	8	7	1		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	B	1	Total	C	F	N	Ru	0	1
			70	54	8	7	1		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	391	Total	O	0	0
			391	391		
6	B	303	Total	O	0	0
			303	303		

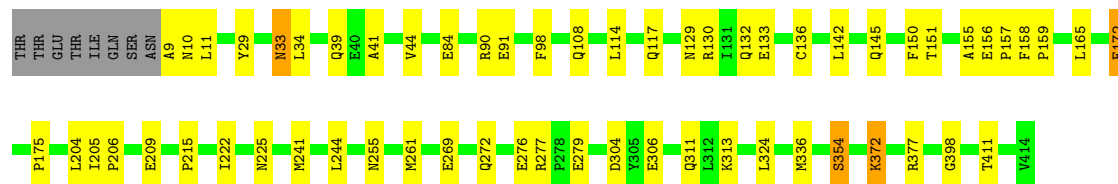
3 Residue-property plots

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. 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. 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.

Note EDS was not executed.

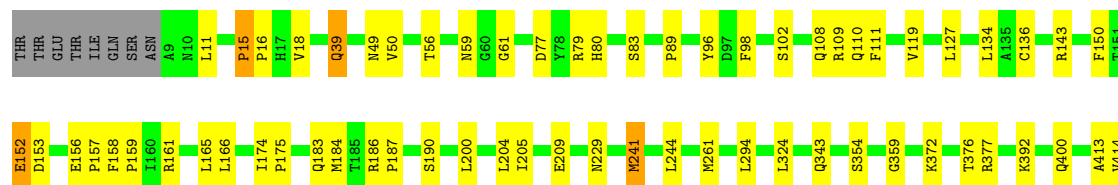
• Molecule 1: Cytochrome P450CAM

Chain A: 



• Molecule 1: Cytochrome P450CAM

Chain B: 



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	63.87Å 67.05Å 72.52Å 71.16° 65.20° 62.31°	Depositor
Resolution (Å)	19.81 – 1.65	Depositor
% Data completeness (in resolution range)	97.8 (19.81-1.65)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	0.04	Depositor
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.210 , 0.226	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	7649	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: CAC, RFB, RFA, 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	0.45	1/3390 (0.0%)	0.90	4/4605 (0.1%)
1	B	0.42	0/3337	0.90	10/4532 (0.2%)
All	All	0.43	1/6727 (0.0%)	0.90	14/9137 (0.2%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	336	MET	SD-CE	-6.58	1.63	1.79

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	324	LEU	N-CA-C	7.91	121.00	111.82
1	B	109	ARG	N-CA-C	6.81	119.28	111.11
1	B	50	VAL	N-CA-C	6.51	114.98	107.89
1	A	204	LEU	N-CA-C	6.13	117.76	111.14
1	B	324	LEU	N-CA-C	6.03	118.82	111.82
1	B	18	VAL	N-CA-C	5.78	114.24	107.77
1	B	15	PRO	CA-C-N	5.60	125.70	119.32
1	B	15	PRO	C-N-CA	5.60	125.70	119.32
1	B	153	ASP	N-CA-C	5.48	119.69	112.89
1	B	204	LEU	N-CA-C	5.41	116.98	111.14
1	B	186	ARG	CA-C-N	5.31	125.29	120.03
1	B	186	ARG	C-N-CA	5.31	125.29	120.03
1	A	34	LEU	N-CA-C	-5.25	106.13	112.54
1	A	354	SER	N-CA-C	5.19	117.67	111.71

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	3311	0	3243	45	0
1	B	3258	0	3196	43	0
2	A	8	0	0	1	0
2	B	12	0	0	5	0
3	A	43	0	30	0	0
3	B	43	0	30	2	0
4	A	70	0	60	0	0
4	B	70	0	60	1	0
5	A	70	0	60	1	0
5	B	70	0	60	1	0
6	A	391	0	0	8	0
6	B	303	0	0	3	0
All	All	7649	0	6739	88	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 6.

All (88) 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:110:GLN:HE22	1:B:229:ASN:H	1.13	0.96
1:B:136[A]:CYS:HB3	2:B:504:CAC:AS	2.30	0.92
1:A:306:GLU:HG3	1:A:311:GLN:HE22	1.38	0.89
1:A:306:GLU:CG	1:A:311:GLN:HE22	1.93	0.81
1:A:306:GLU:HG3	1:A:311:GLN:NE2	1.96	0.79
1:A:151[A]:THR:HG22	6:A:1278:HOH:O	1.85	0.76
1:A:306:GLU:CD	1:A:311:GLN:HE22	1.94	0.76
1:A:132:GLN:NE2	1:A:136:CYS:SG	2.61	0.73
1:A:11:LEU:HD23	2:A:503:CAC:C1	2.21	0.70
1:B:158:PHE:HB3	1:B:159:PRO:HD3	1.75	0.69
1:B:136[A]:CYS:CB	2:B:504:CAC:AS	2.98	0.68
1:A:130:ARG:HG2	1:A:130:ARG:HH21	1.59	0.66
1:B:376:THR:CG2	1:B:414:VAL:HG21	2.29	0.63
1:A:241[A]:MET:HG3	6:A:1562:HOH:O	1.99	0.62
1:A:158:PHE:HB3	1:A:159:PRO:HD3	1.80	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:84:GLU:CD	1:A:90:ARG:HH12	2.08	0.61
1:B:110:GLN:NE2	1:B:229:ASN:H	1.91	0.60
1:B:108:GLN:HE22	1:B:354:SER:HB2	1.67	0.60
1:A:255[A]:ASN:OD1	1:A:398:GLY:HA3	2.02	0.59
1:A:313:LYS:HE3	6:A:1617:HOH:O	2.03	0.59
1:B:11:LEU:HD23	2:B:502:CAC:C2	2.33	0.59
1:A:304:ASP:OD1	1:A:313:LYS:HD2	2.04	0.58
1:B:343[A]:GLN:HA	1:B:343[A]:GLN:OE1	2.02	0.58
1:B:119:VAL:HG12	1:B:166:LEU:HD21	1.85	0.58
1:B:136[B]:CYS:SG	1:B:377:ARG:NH2	2.77	0.58
1:A:29:TYR:CD2	5:A:901[B]:RFB:H61	2.40	0.57
1:B:127[A]:LEU:HD11	1:B:165:LEU:HG	1.85	0.57
3:B:417:HEM:HBC2	3:B:417:HEM:HHD	1.86	0.56
1:A:108:GLN:HE22	1:A:354:SER:HB2	1.71	0.55
1:B:134:LEU:HD21	1:B:161:ARG:HB2	1.89	0.55
1:B:136[A]:CYS:HB2	2:B:504:CAC:C2	2.37	0.55
1:A:9:ALA:O	1:A:10:ASN:HB2	2.06	0.55
1:A:277:ARG:HD3	1:A:279:GLU:OE1	2.07	0.54
1:B:372:LYS:HD3	6:B:1359:HOH:O	2.07	0.54
1:A:306:GLU:HG2	6:A:1466:HOH:O	2.06	0.54
1:B:376:THR:HG22	1:B:414:VAL:HG21	1.89	0.54
1:B:96:TYR:HB3	2:B:500:CAC:C1	2.39	0.53
1:A:98:PHE:HB3	1:A:244:LEU:HB2	1.91	0.53
1:A:130:ARG:HG2	1:A:130:ARG:NH2	2.22	0.53
1:A:172:GLU:H	1:A:172:GLU:CD	2.17	0.53
1:A:44[A]:VAL:HG11	6:A:1503:HOH:O	2.09	0.53
1:A:175:PRO:HG3	1:B:79:ARG:HH11	1.74	0.52
1:A:269[B]:GLU:HG2	6:A:1424:HOH:O	2.10	0.52
1:B:156:GLU:HB2	1:B:157:PRO:HD3	1.91	0.52
1:A:130:ARG:NH2	1:A:165:LEU:HD11	2.26	0.51
1:B:111:PHE:HD2	1:B:241[B]:MET:SD	2.34	0.50
1:B:134:LEU:HD11	1:B:161:ARG:HD2	1.92	0.50
1:B:205:ILE:O	1:B:209[B]:GLU:HG2	2.12	0.50
1:A:41:ALA:O	1:A:44[A]:VAL:HG22	2.11	0.49
1:B:187:PRO:HG3	4:B:902[A]:RFA:H211	1.95	0.48
1:A:129:ASN:O	1:A:133:GLU:HG3	2.13	0.48
1:A:205:ILE:O	1:A:209:GLU:HG2	2.13	0.48
1:A:372:LYS:HE3	1:A:372:LYS:HB2	1.66	0.48
1:B:77:ASP:OD2	1:B:80:HIS:HB2	2.14	0.48
1:B:98:PHE:HB3	1:B:244:LEU:HB2	1.96	0.47
1:A:272[B]:GLN:HE21	1:A:276[B]:GLU:CD	2.23	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:184:MET:HE1	1:B:200:LEU:HD22	1.98	0.46
1:A:311:GLN:NE2	1:A:311:GLN:HA	2.31	0.46
1:A:33[A]:ASN:CG	1:A:33[A]:ASN:O	2.58	0.45
1:A:215:PRO:HB3	1:A:225:ASN:OD1	2.15	0.45
1:B:187:PRO:HG3	5:B:903[B]:RFB:H211	1.98	0.45
1:B:56:THR:O	1:B:61:GLY:HA2	2.17	0.44
1:B:392:LYS:HD3	1:B:400:GLN:OE1	2.17	0.44
1:A:150:PHE:CZ	1:A:261:MET:HG3	2.52	0.44
1:B:59:ASN:HB3	1:B:89:PRO:HD3	1.99	0.44
1:A:142:LEU:HA	1:A:145:GLN:NE2	2.33	0.44
1:A:151[A]:THR:HA	1:A:155:ALA:HB3	2.00	0.44
1:B:294:LEU:H	1:B:294:LEU:HD23	1.83	0.43
1:A:151[B]:THR:HA	1:A:155:ALA:HB3	2.01	0.43
1:A:205:ILE:HB	1:A:206:PRO:HD3	2.01	0.42
1:B:39:GLN:NE2	1:B:39:GLN:H	2.16	0.42
1:A:151[B]:THR:HG21	6:A:1594:HOH:O	2.19	0.42
1:A:91:GLU:H	1:A:91:GLU:CD	2.28	0.42
1:B:143:ARG:NH2	1:B:413:ALA:HB2	2.35	0.42
1:B:241[A]:MET:HG3	6:B:1203:HOH:O	2.20	0.42
1:B:83[A]:SER:O	1:B:102:SER:HA	2.20	0.41
1:B:152:GLU:HG2	6:B:1396:HOH:O	2.20	0.41
1:B:134:LEU:HD11	1:B:161:ARG:CD	2.49	0.41
1:B:183:GLN:OE1	1:B:190:SER:HB2	2.20	0.41
1:B:15:PRO:HA	1:B:16:PRO:HD3	1.86	0.41
1:B:174:ILE:HB	1:B:175:PRO:HD3	2.03	0.41
1:A:130:ARG:HD3	6:A:1273:HOH:O	2.21	0.41
1:B:150:PHE:CZ	1:B:261:MET:HG3	2.55	0.40
1:B:359:GLY:HA3	3:B:417:HEM:C3C	2.56	0.40
1:A:156:GLU:HB2	1:A:157:PRO:HD3	2.04	0.40
1:B:83[B]:SER:O	1:B:102:SER:HA	2.20	0.40
1:A:114:LEU:O	1:A:117:GLN:HB2	2.21	0.40
1:A:377:ARG:O	1:A:411:THR:HB	2.22	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	416/414 (100%)	398 (96%)	18 (4%)	0	100	100
1	B	410/414 (99%)	396 (97%)	14 (3%)	0	100	100
All	All	826/828 (100%)	794 (96%)	32 (4%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	361/357 (101%)	355 (98%)	6 (2%)	53	32
1	B	355/357 (99%)	350 (99%)	5 (1%)	59	39
All	All	716/714 (100%)	705 (98%)	11 (2%)	61	37

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	33[A]	ASN
1	A	33[B]	ASN
1	A	39	GLN
1	A	172	GLU
1	A	222	ILE
1	A	372	LYS
1	B	39	GLN
1	B	49	ASN
1	B	152	GLU
1	B	241[A]	MET
1	B	241[B]	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	39	GLN
1	A	46	GLN
1	A	69	GLN
1	A	108	GLN
1	A	132	GLN
1	A	147	GLN
1	A	210	GLN
1	A	311	GLN
1	B	17	HIS
1	B	21	HIS
1	B	30	ASN
1	B	33	ASN
1	B	39	GLN
1	B	46	GLN
1	B	108	GLN
1	B	110	GLN
1	B	116	ASN
1	B	117	GLN
1	B	147	GLN
1	B	210	GLN
1	B	227	GLN
1	B	311	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](#)

11 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	HEM	A	417	1,6	50,50,50	1.36	6 (12%)	67,82,82	0.86	2 (2%)
4	RFA	A	900[A]	-	77,83,83	1.48	11 (14%)	87,137,137	1.62	12 (13%)
2	CAC	B	500	1	0,3,4	-	-	0,3,6	-	-
2	CAC	B	504	1	0,3,4	-	-	0,3,6	-	-
5	RFB	A	901[B]	-	77,83,83	1.53	13 (16%)	87,137,137	2.01	15 (17%)
2	CAC	A	501	1	0,3,4	-	-	0,3,6	-	-
2	CAC	A	503	1	0,3,4	-	-	0,3,6	-	-
4	RFA	B	902[A]	-	77,83,83	1.51	10 (12%)	87,137,137	1.75	15 (17%)
5	RFB	B	903[B]	-	77,83,83	1.57	14 (18%)	87,137,137	1.99	14 (16%)
2	CAC	B	502	1	0,3,4	-	-	0,3,6	-	-
3	HEM	B	417	1,6	50,50,50	1.35	6 (12%)	67,82,82	0.95	3 (4%)

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	A	417	1,6	-	2/14/54/54	-
4	RFA	A	900[A]	-	-	0/12/182/182	0/15/14/14
5	RFB	A	901[B]	-	-	1/12/182/182	0/15/14/14
4	RFA	B	902[A]	-	-	2/12/182/182	0/15/14/14
5	RFB	B	903[B]	-	-	1/12/182/182	0/15/14/14
3	HEM	B	417	1,6	-	3/14/54/54	-

All (60) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	902[A]	RFA	C50-C43	-5.37	1.39	1.50
4	A	900[A]	RFA	C50-C43	-4.66	1.40	1.50
5	A	901[B]	RFB	C50-C43	-4.59	1.40	1.50
5	B	903[B]	RFB	C50-C43	-3.84	1.42	1.50
5	B	903[B]	RFB	C43-C42	3.37	1.44	1.39
3	A	417	HEM	FE-ND	3.31	2.05	1.94

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	417	HEM	FE-ND	3.31	2.05	1.94
3	B	417	HEM	FE-NA	3.25	2.05	1.95
3	A	417	HEM	FE-NA	3.16	2.05	1.95
4	B	902[A]	RFA	C43-C42	3.13	1.44	1.39
3	B	417	HEM	FE-NB	3.12	2.04	1.94
5	B	903[B]	RFB	C63-C61	3.04	1.61	1.53
4	B	902[A]	RFA	C62-C63	3.02	1.60	1.53
3	B	417	HEM	FE-NC	2.99	2.05	1.95
5	B	903[B]	RFB	C50-C55	2.97	1.43	1.39
4	A	900[A]	RFA	C63-C61	2.85	1.60	1.53
5	A	901[B]	RFB	C50-C55	2.83	1.43	1.39
4	B	902[A]	RFA	F59-C55	2.82	1.39	1.34
4	B	902[A]	RFA	C50-C51	2.77	1.43	1.39
4	A	900[A]	RFA	C43-C42	2.76	1.43	1.39
4	A	900[A]	RFA	C50-C55	2.76	1.43	1.39
5	B	903[B]	RFB	C43-C44	2.74	1.43	1.39
5	B	903[B]	RFB	C50-C51	2.68	1.43	1.39
3	A	417	HEM	FE-NB	2.66	2.03	1.94
3	B	417	HEM	CAB-C3B	-2.64	1.40	1.47
5	A	901[B]	RFB	C63-C61	2.64	1.60	1.53
5	A	901[B]	RFB	C66-C61	2.64	1.60	1.53
3	A	417	HEM	FE-NC	2.63	2.03	1.95
3	B	417	HEM	CAC-C3C	-2.60	1.40	1.47
3	A	417	HEM	CAC-C3C	-2.49	1.40	1.47
3	A	417	HEM	CAB-C3B	-2.46	1.40	1.47
5	A	901[B]	RFB	C43-C42	2.46	1.42	1.39
4	B	902[A]	RFA	C52-C51	2.43	1.41	1.38
4	A	900[A]	RFA	C50-C51	2.42	1.42	1.39
5	A	901[B]	RFB	C68-C66	2.39	1.58	1.53
5	A	901[B]	RFB	C55-C54	2.38	1.41	1.38
4	B	902[A]	RFA	C62-C69	2.35	1.59	1.52
5	A	901[B]	RFB	C62-C63	2.32	1.58	1.53
4	B	902[A]	RFA	C68-C69	2.32	1.59	1.52
5	A	901[B]	RFB	C50-C51	2.32	1.42	1.39
5	B	903[B]	RFB	C68-C66	2.32	1.58	1.53
4	A	900[A]	RFA	C68-C66	2.24	1.58	1.53
5	B	903[B]	RFB	C55-C54	2.22	1.41	1.38
5	B	903[B]	RFB	C67-C66	2.21	1.58	1.53
5	B	903[B]	RFB	C62-C63	2.20	1.58	1.53
5	B	903[B]	RFB	C20-N25	2.20	1.38	1.34
5	A	901[B]	RFB	C53-C54	2.17	1.42	1.39
5	B	903[B]	RFB	C66-C61	2.14	1.59	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	900[A]	RFA	C62-C63	2.12	1.58	1.53
4	A	900[A]	RFA	C67-C66	2.11	1.58	1.53
5	B	903[B]	RFB	F47-C42	2.10	1.37	1.34
4	A	900[A]	RFA	C43-C44	2.08	1.42	1.39
4	B	902[A]	RFA	C50-C55	2.08	1.42	1.39
5	A	901[B]	RFB	C39-C40	-2.06	1.47	1.51
4	A	900[A]	RFA	C66-C61	2.06	1.58	1.53
4	B	902[A]	RFA	C67-C65	2.05	1.58	1.52
5	B	903[B]	RFB	C19-N14	2.05	1.37	1.34
5	A	901[B]	RFB	C62-C69	2.04	1.58	1.52
4	A	900[A]	RFA	C62-C69	2.02	1.58	1.52
5	A	901[B]	RFB	C43-C44	2.01	1.42	1.39

All (61) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	901[B]	RFB	C18-C19-N14	-8.97	115.29	124.84
5	B	903[B]	RFB	C18-C19-N14	-8.70	115.57	124.84
5	A	901[B]	RFB	C21-C20-N25	-7.46	116.89	124.84
5	B	903[B]	RFB	C21-C20-N25	-7.46	116.90	124.84
4	A	900[A]	RFA	C18-C19-N14	-6.07	118.37	124.84
5	B	903[B]	RFB	C6-C7-N2	-5.88	118.58	124.84
5	B	903[B]	RFB	C30-C31-N26	-5.77	118.69	124.84
5	A	901[B]	RFB	C30-C31-N26	-5.74	118.73	124.84
5	A	901[B]	RFB	C6-C7-N2	-5.63	118.85	124.84
4	B	902[A]	RFA	C18-C19-N14	-5.60	118.88	124.84
4	B	902[A]	RFA	C30-C31-N26	-5.23	119.27	124.84
4	B	902[A]	RFA	C6-C7-N2	-5.23	119.27	124.84
4	A	900[A]	RFA	C6-C7-N2	-5.11	119.40	124.84
4	A	900[A]	RFA	C30-C31-N26	-4.92	119.60	124.84
4	A	900[A]	RFA	C21-C20-N25	-4.47	120.08	124.84
4	B	902[A]	RFA	C21-C20-N25	-4.37	120.19	124.84
4	A	900[A]	RFA	C9-C8-N13	-4.30	120.26	124.84
4	B	902[A]	RFA	C9-C8-N13	-4.28	120.28	124.84
4	B	902[A]	RFA	C45-C40-C41	4.12	120.75	115.91
4	B	902[A]	RFA	C33-C32-N37	-4.02	120.56	124.84
5	B	903[B]	RFB	C9-C8-N13	-4.00	120.59	124.84
4	A	900[A]	RFA	C33-C32-N37	-3.95	120.64	124.84
4	A	900[A]	RFA	C45-C40-C41	3.78	120.35	115.91
5	B	903[B]	RFB	C45-C40-C41	3.72	120.28	115.91
5	A	901[B]	RFB	C39-C40-C41	-3.69	118.12	122.13
5	A	901[B]	RFB	C45-C40-C41	3.68	120.23	115.91

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	903[B]	RFB	C39-C40-C41	-3.65	118.16	122.13
5	A	901[B]	RFB	C38-C17-C16	3.64	127.68	120.00
5	B	903[B]	RFB	C38-C17-C16	3.61	127.61	120.00
5	A	901[B]	RFB	C9-C8-N13	-3.60	121.01	124.84
4	B	902[A]	RFA	C54-C53-C52	3.37	121.87	115.99
3	B	417	HEM	CAC-C3C-C4C	3.28	132.65	124.82
5	A	901[B]	RFB	C33-C32-N37	-3.24	121.39	124.84
5	B	903[B]	RFB	C33-C32-N37	-3.07	121.57	124.84
3	A	417	HEM	C4B-C3B-C2B	-3.01	104.51	107.28
3	A	417	HEM	C3B-C4B-NB	2.89	111.54	109.47
4	B	902[A]	RFA	C39-C22-C21	-2.76	120.04	123.33
4	B	902[A]	RFA	C66-C61-N60	2.66	117.32	112.00
5	A	901[B]	RFB	C27-C28-C29	-2.49	117.13	120.51
5	A	901[B]	RFB	C38-C17-C18	-2.48	116.33	121.23
4	B	902[A]	RFA	C39-C40-C41	-2.41	119.51	122.13
5	B	903[B]	RFB	C27-C28-C29	-2.41	117.24	120.51
3	B	417	HEM	CAC-C3C-C2C	-2.40	120.62	128.43
5	B	903[B]	RFB	C3-C4-C5	-2.36	117.31	120.51
5	A	901[B]	RFB	C39-C22-C21	-2.33	120.54	123.33
5	B	903[B]	RFB	C38-C17-C18	-2.32	116.64	121.23
5	A	901[B]	RFB	C3-C4-C5	-2.27	117.43	120.51
4	A	900[A]	RFA	C39-C22-C21	-2.26	120.64	123.33
5	B	903[B]	RFB	C39-C22-C21	-2.23	120.67	123.33
4	B	902[A]	RFA	C40-C45-C44	-2.21	119.83	122.36
4	B	902[A]	RFA	C39-C40-C45	-2.20	119.74	122.13
4	B	902[A]	RFA	C53-N60-C61	2.20	133.18	121.08
3	B	417	HEM	C3B-C4B-NB	2.19	111.04	109.47
4	A	900[A]	RFA	C39-C40-C45	-2.18	119.76	122.13
5	B	903[B]	RFB	C20-C21-C22	2.18	122.59	120.28
5	A	901[B]	RFB	C20-C21-C22	2.14	122.55	120.28
4	B	902[A]	RFA	C27-C28-C29	-2.11	117.65	120.51
4	A	900[A]	RFA	C27-C28-C29	-2.07	117.70	120.51
4	A	900[A]	RFA	C39-C40-C41	-2.06	119.89	122.13
4	A	900[A]	RFA	C40-C41-C42	-2.02	120.05	122.36
5	A	901[B]	RFB	C15-C16-C17	-2.01	116.72	120.99

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	902[A]	RFA	C63-C61-N60-C53
5	B	903[B]	RFB	C63-C61-N60-C53

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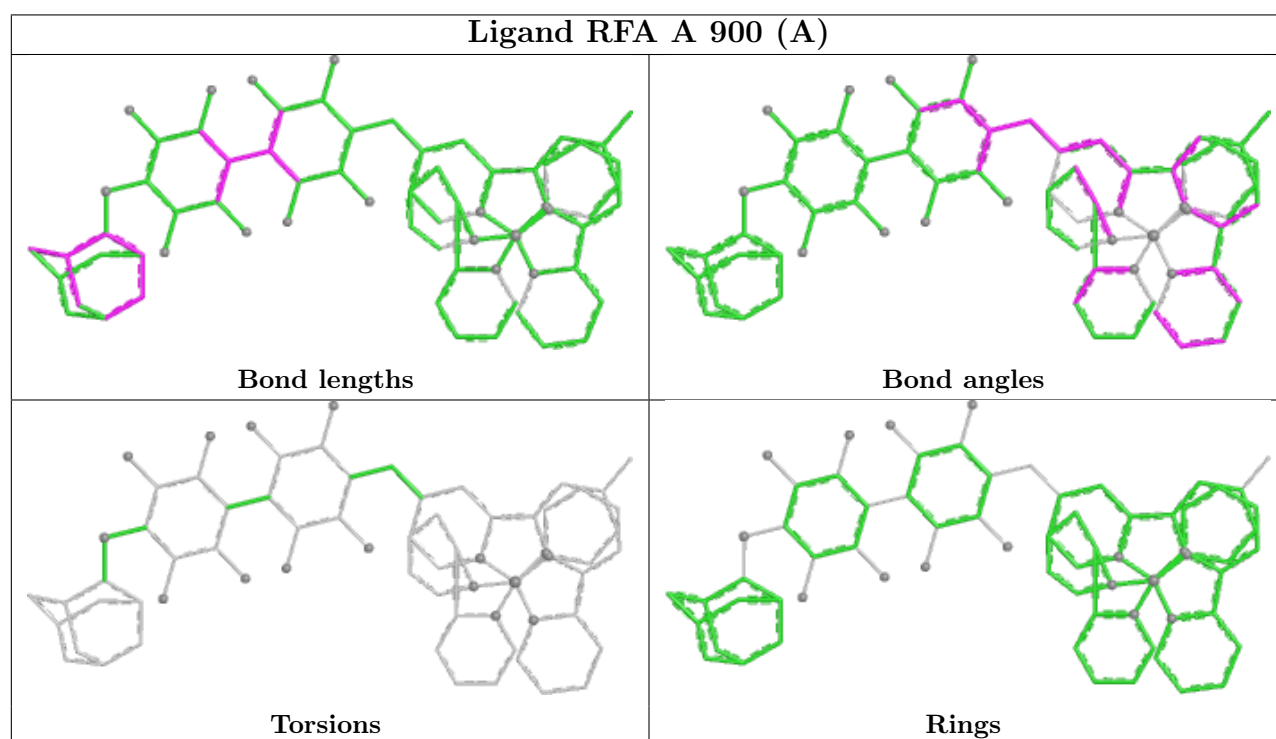
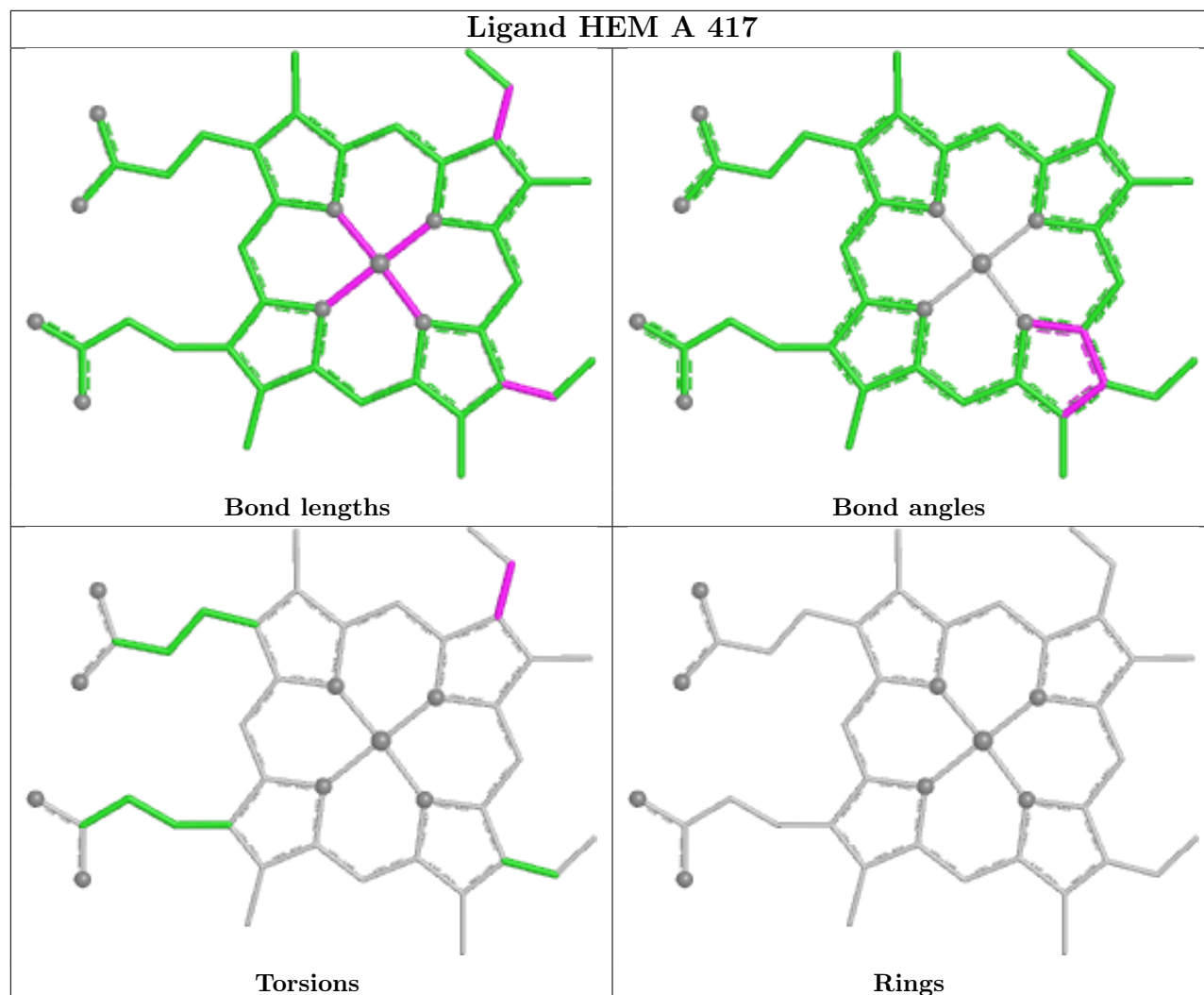
Mol	Chain	Res	Type	Atoms
4	B	902[A]	RFA	C21-C22-C39-C40
3	A	417	HEM	C2C-C3C-CAC-CBC
3	B	417	HEM	C4C-C3C-CAC-CBC
3	A	417	HEM	C4C-C3C-CAC-CBC
3	B	417	HEM	CAD-CBD-CGD-O1D
3	B	417	HEM	CAD-CBD-CGD-O2D
5	A	901[B]	RFB	C54-C53-N60-C61

There are no ring outliers.

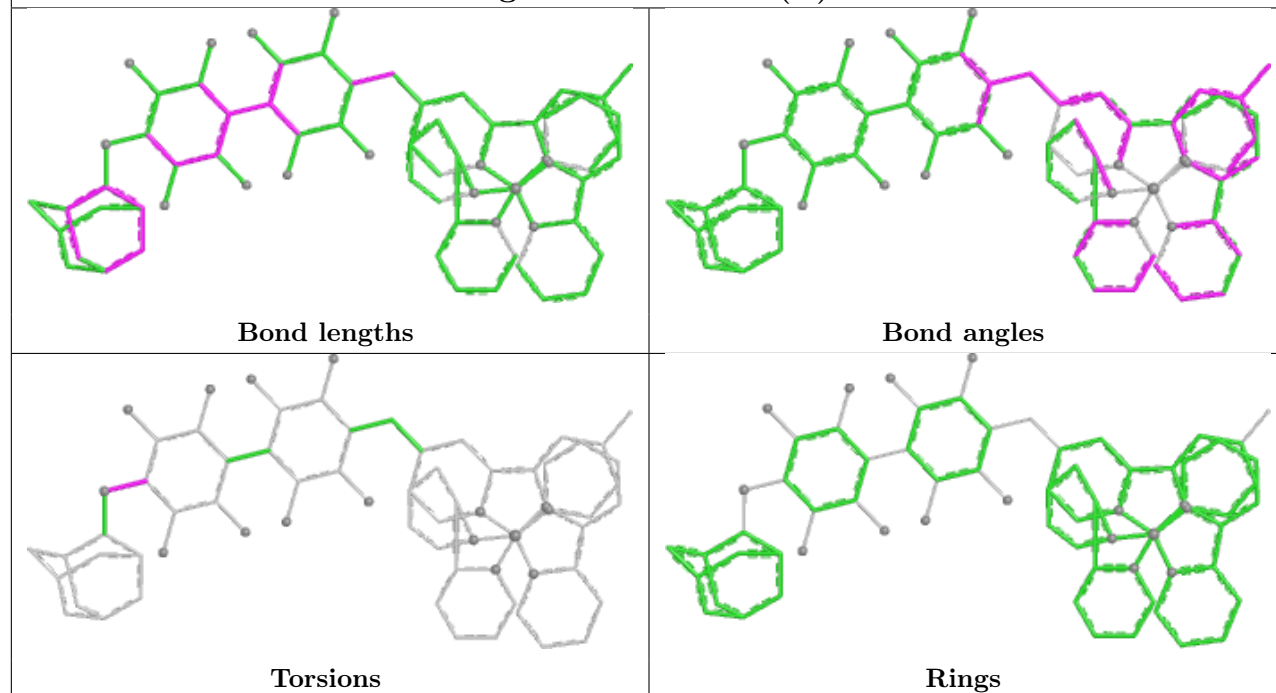
8 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	500	CAC	1	0
2	B	504	CAC	3	0
5	A	901[B]	RFB	1	0
2	A	503	CAC	1	0
4	B	902[A]	RFA	1	0
5	B	903[B]	RFB	1	0
2	B	502	CAC	1	0
3	B	417	HEM	2	0

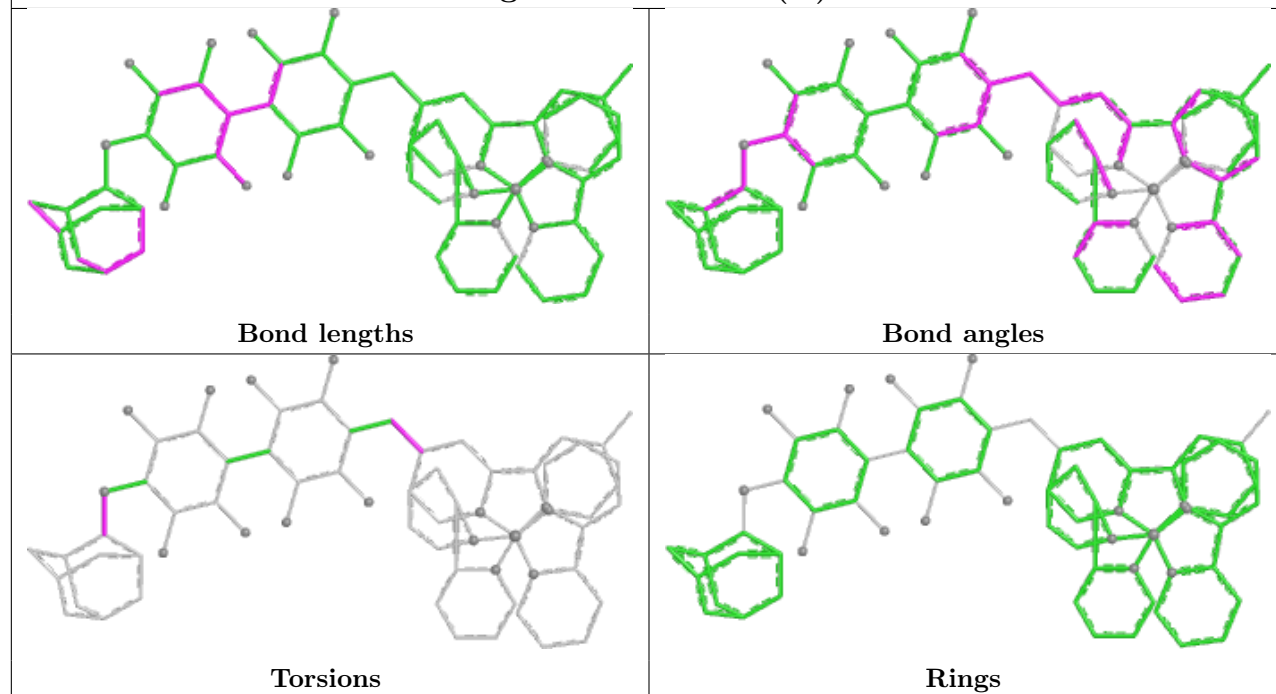
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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



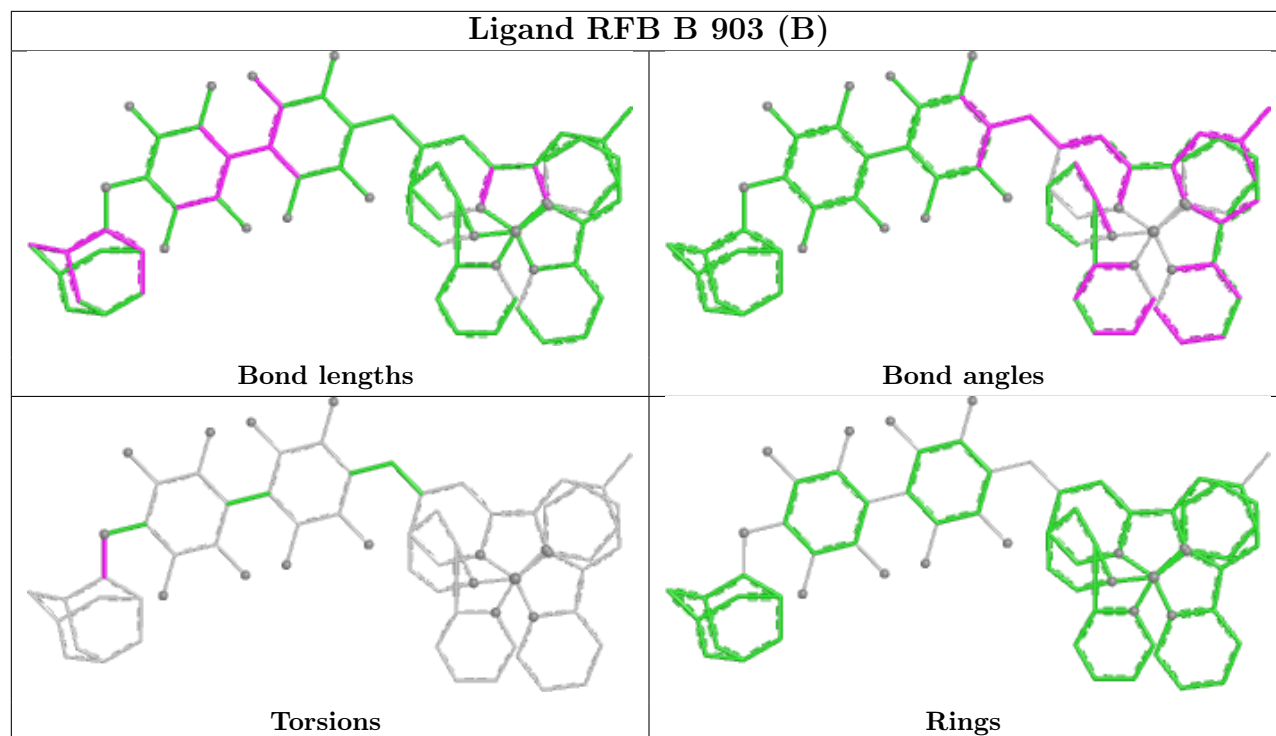
Ligand RFB A 901 (B)



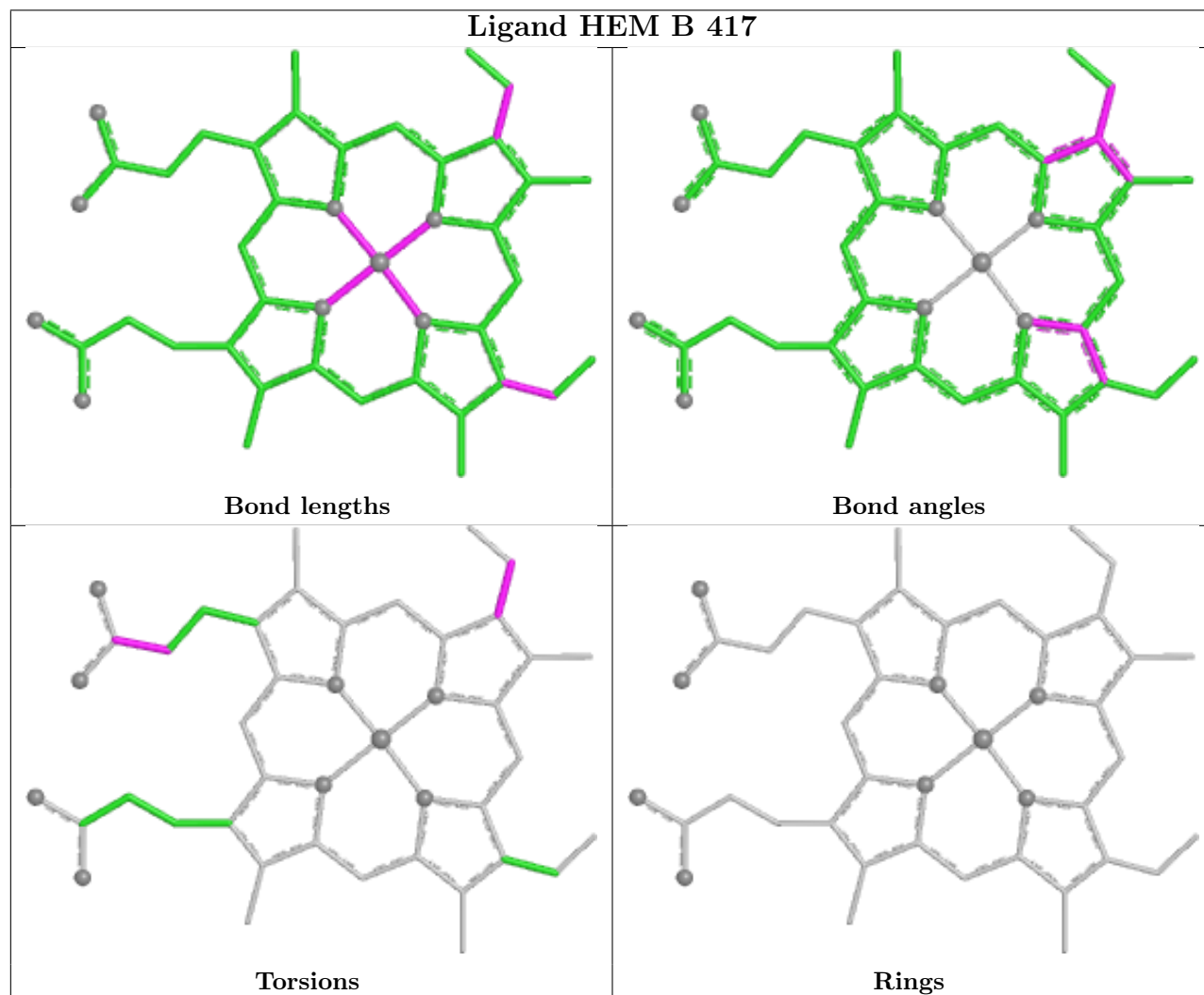
Ligand RFA B 902 (A)



Ligand RFB B 903 (B)



Ligand HEM B 417



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

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.