



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 10:39 PM UTC

PDB ID : 1PHB / pdb_00001phb
Title : INHIBITOR-INDUCED CONFORMATIONAL CHANGE IN CY-TOCHROME P450-CAM
Authors : Poulos, T.L.
Deposited on : 1992-07-27
Resolution : 1.60 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) : 2.49

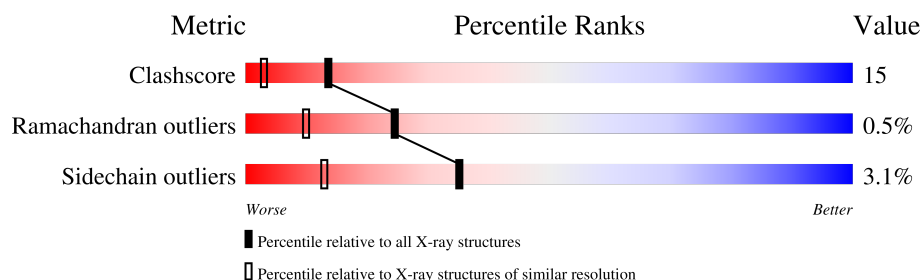
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.60 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	4931 (1.60-1.60)
Ramachandran outliers	187476	4831 (1.60-1.60)
Sidechain outliers	187428	4830 (1.60-1.60)

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	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

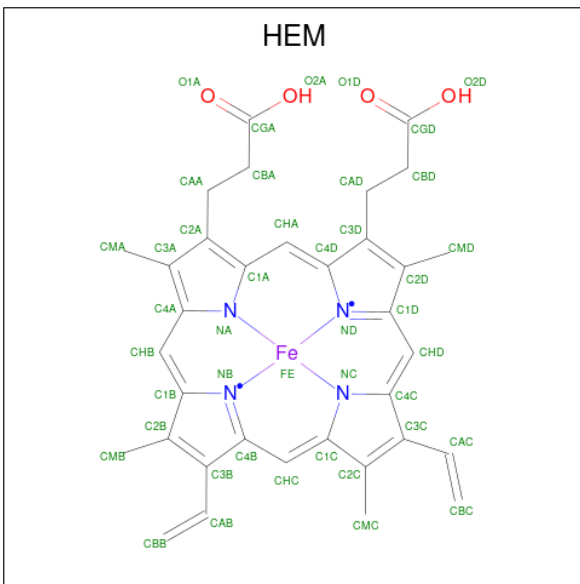
Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	PFZ	A	422	-	-	X	-

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 P450-CAM.

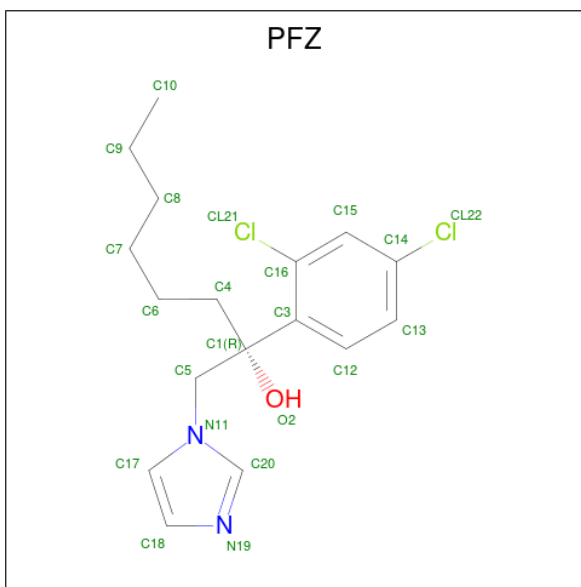
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	405	Total	C	N	O	S	0	4	0
			3227	2046	564	598	19			

- 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 43	C 34	Fe 1	N 4	O 4	0	0

- Molecule 3 is 1-(N-IMIDAZOLYL)-2-HYDROXY-2-(2,3-DICHLOROPHENYL)OCTANE (CCD ID: PFZ) (formula: $C_{17}H_{22}Cl_2N_2O$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	Cl	N	O	0	0
			22	17	2	2	1		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	213	Total	O	0	0
			213	213		

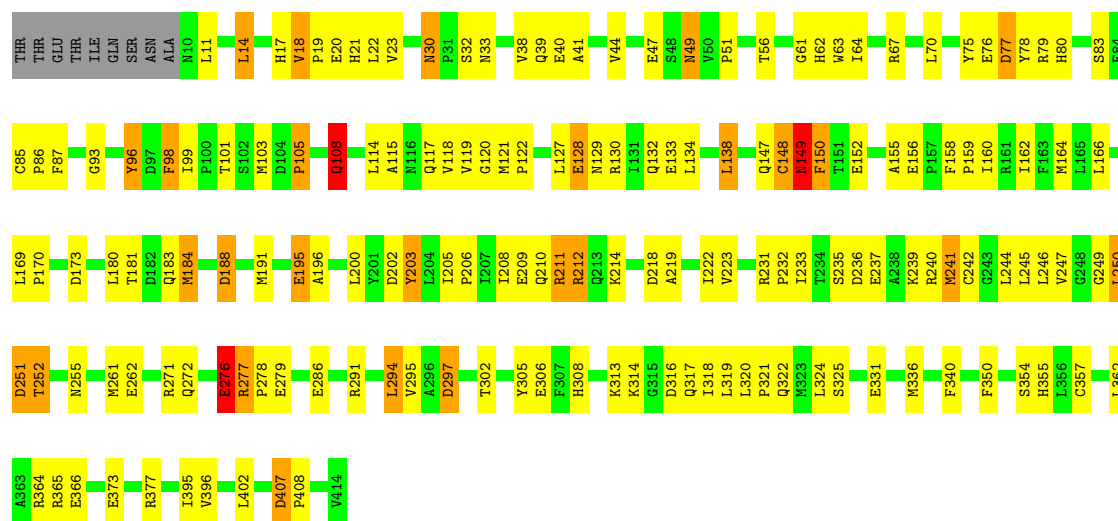
3 Residue-property plots [i](#)

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. 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 P450-CAM

Chain A: 



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	108.67Å 103.90Å 36.38Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	(Not available) – 1.60	Depositor
% Data completeness (in resolution range)	(Not available) ((Not available)-1.60)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	PROLSQ	Depositor
R, R_{free}	0.190 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3505	wwPDB-VP
Average B, all atoms (Å ²)	20.0	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: PFZ, 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	1.33	4/3321 (0.1%)	2.07	110/4510 (2.4%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	1	8

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	211	ARG	CD-NE	-6.12	1.37	1.46
1	A	149	ASN	N-CA	-5.95	1.39	1.46
1	A	85	CYS	N-CA	5.54	1.52	1.46
1	A	396	VAL	N-CA	5.07	1.51	1.46

All (110) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	211	ARG	CD-NE-CZ	18.04	149.65	124.40
1	A	148	CYS	CA-C-N	13.62	142.12	122.24
1	A	148	CYS	C-N-CA	13.62	142.12	122.24
1	A	149	ASN	N-CA-CB	10.07	121.94	110.35
1	A	33	ASN	CA-CB-CG	9.02	121.62	112.60
1	A	32	SER	N-CA-C	8.96	120.81	111.14
1	A	316	ASP	CA-CB-CG	8.59	121.19	112.60
1	A	96	TYR	N-CA-CB	8.49	124.17	110.16
1	A	33	ASN	CB-CA-C	8.47	123.35	111.74

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	250	LEU	N-CA-C	8.31	121.91	111.69
1	A	96	TYR	CB-CA-C	8.08	123.05	109.80
1	A	286	GLU	N-CA-CB	7.97	121.96	110.16
1	A	324	LEU	N-CA-C	7.36	119.38	111.36
1	A	188	ASP	N-CA-CB	-7.32	99.98	110.81
1	A	252	THR	CA-C-N	7.28	130.44	120.46
1	A	252	THR	C-N-CA	7.28	130.44	120.46
1	A	149	ASN	CA-C-O	-6.96	111.28	122.20
1	A	202	ASP	CA-C-N	6.91	129.78	120.65
1	A	202	ASP	C-N-CA	6.91	129.78	120.65
1	A	291	ARG	NE-CZ-NH2	6.86	125.38	119.20
1	A	407	ASP	CB-CA-C	6.81	117.61	109.85
1	A	180	LEU	O-C-N	6.76	129.29	122.12
1	A	30	ASN	O-C-N	6.75	129.08	121.32
1	A	33	ASN	N-CA-CB	6.74	118.10	110.35
1	A	32	SER	CA-C-O	-6.71	113.79	120.70
1	A	251	ASP	CB-CA-C	6.69	121.41	110.24
1	A	98	PHE	CA-CB-CG	6.65	120.45	113.80
1	A	20	GLU	CA-CB-CG	6.63	127.36	114.10
1	A	64	ILE	CA-C-O	-6.61	113.70	120.25
1	A	118	VAL	N-CA-C	6.58	120.26	111.44
1	A	279	GLU	CB-CG-CD	6.43	123.54	112.60
1	A	119	VAL	CB-CA-C	6.24	116.77	110.65
1	A	80	HIS	CA-CB-CG	-6.18	107.62	113.80
1	A	21	HIS	CA-CB-CG	6.15	119.95	113.80
1	A	152	GLU	N-CA-CB	6.11	119.30	110.20
1	A	211	ARG	CG-CD-NE	6.10	125.43	112.00
1	A	62	HIS	CA-CB-CG	6.07	119.87	113.80
1	A	76	GLU	CA-C-N	6.06	133.11	121.54
1	A	76	GLU	C-N-CA	6.06	133.11	121.54
1	A	209	GLU	CA-C-N	6.03	128.68	120.54
1	A	209	GLU	C-N-CA	6.03	128.68	120.54
1	A	108	GLN	CA-CB-CG	5.88	125.85	114.10
1	A	241	MET	N-CA-CB	5.76	118.43	110.07
1	A	149	ASN	CB-CA-C	5.73	119.59	111.74
1	A	44	VAL	CA-C-N	5.73	129.77	120.60
1	A	44	VAL	C-N-CA	5.73	129.77	120.60
1	A	67	ARG	NE-CZ-NH2	-5.71	114.06	119.20
1	A	252	THR	N-CA-C	5.68	122.90	110.80
1	A	210	GLN	CA-CB-CG	5.64	125.38	114.10
1	A	130	ARG	CA-CB-CG	5.64	125.38	114.10
1	A	18	VAL	N-CA-C	5.61	112.66	107.56

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	148	CYS	O-C-N	-5.59	113.83	122.61
1	A	156	GLU	CB-CG-CD	5.58	122.09	112.60
1	A	297	ASP	N-CA-CB	5.57	119.90	110.49
1	A	164	MET	CA-C-N	5.56	128.19	120.29
1	A	164	MET	C-N-CA	5.56	128.19	120.29
1	A	245	LEU	CA-C-N	5.54	127.65	120.44
1	A	245	LEU	C-N-CA	5.54	127.65	120.44
1	A	152	GLU	CA-CB-CG	5.54	125.19	114.10
1	A	262	GLU	CB-CG-CD	5.52	121.99	112.60
1	A	133	GLU	CA-C-O	-5.51	115.03	120.82
1	A	203	TYR	CA-C-N	5.51	127.61	120.44
1	A	203	TYR	C-N-CA	5.51	127.61	120.44
1	A	105	PRO	N-CA-C	-5.50	105.19	110.47
1	A	138	LEU	CB-CA-C	5.50	119.59	110.90
1	A	188	ASP	CA-CB-CG	5.47	118.07	112.60
1	A	239	LYS	CA-C-N	5.46	127.59	120.28
1	A	239	LYS	C-N-CA	5.46	127.59	120.28
1	A	169	LEU	O-C-N	5.45	126.86	121.57
1	A	14	LEU	O-C-N	5.45	126.15	121.30
1	A	77	ASP	N-CA-C	5.45	122.40	110.80
1	A	340	PHE	CA-CB-CG	5.43	119.23	113.80
1	A	51	PRO	CA-C-N	5.43	127.56	120.28
1	A	51	PRO	C-N-CA	5.43	127.56	120.28
1	A	129	ASN	CA-C-O	-5.43	114.80	120.55
1	A	11	LEU	CB-CA-C	5.42	118.58	109.48
1	A	21	HIS	N-CA-CB	5.42	119.23	110.40
1	A	130	ARG	N-CA-CB	5.40	118.16	110.16
1	A	75	TYR	N-CA-CB	5.39	118.64	110.28
1	A	160	ILE	N-CA-CB	5.37	116.83	110.55
1	A	249	GLY	CA-C-N	5.35	129.73	120.58
1	A	249	GLY	C-N-CA	5.35	129.73	120.58
1	A	47	GLU	CB-CG-CD	5.34	121.68	112.60
1	A	39	GLN	N-CA-CB	5.34	117.97	110.12
1	A	38	VAL	CA-C-N	5.33	127.42	120.28
1	A	38	VAL	C-N-CA	5.33	127.42	120.28
1	A	41	ALA	CA-C-N	5.31	127.83	120.29
1	A	41	ALA	C-N-CA	5.31	127.83	120.29
1	A	320	LEU	CA-C-N	5.30	126.47	119.84
1	A	320	LEU	C-N-CA	5.30	126.47	119.84
1	A	322	GLN	OE1-CD-NE2	-5.30	117.30	122.60
1	A	276	GLU	N-CA-C	5.28	118.89	112.23
1	A	115	ALA	CA-C-N	5.27	128.32	120.31

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	115	ALA	C-N-CA	5.27	128.32	120.31
1	A	108	GLN	CB-CG-CD	5.25	121.53	112.60
1	A	128	GLU	N-CA-CB	5.25	117.62	110.01
1	A	306	GLU	CB-CG-CD	5.23	121.49	112.60
1	A	297	ASP	CA-CB-CG	5.22	117.82	112.60
1	A	316	ASP	N-CA-C	-5.20	102.59	110.28
1	A	305	TYR	O-C-N	5.17	129.37	123.48
1	A	138	LEU	N-CA-CB	-5.15	102.60	110.07
1	A	108	GLN	CA-C-N	5.09	127.35	120.38
1	A	108	GLN	C-N-CA	5.09	127.35	120.38
1	A	271	ARG	NE-CZ-NH1	5.08	126.58	121.50
1	A	322	GLN	CB-CG-CD	5.05	121.18	112.60
1	A	23	VAL	N-CA-CB	5.04	116.03	110.53
1	A	294	LEU	N-CA-C	5.04	121.15	113.19
1	A	396	VAL	CA-C-O	-5.03	114.78	120.67
1	A	364	ARG	CD-NE-CZ	5.01	131.42	124.40
1	A	93	GLY	N-CA-C	-5.00	106.69	113.24

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	A	96	TYR	CA

All (8) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	147	GLN	Mainchain
1	A	148	CYS	Mainchain
1	A	149	ASN	Mainchain
1	A	212	ARG	Sidechain
1	A	250	LEU	Mainchain
1	A	277	ARG	Sidechain
1	A	377	ARG	Sidechain
1	A	77	ASP	Mainchain

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	3227	0	3178	85	0
2	A	43	0	30	4	0
3	A	22	0	22	26	0
4	A	213	0	0	9	0
All	All	3505	0	3230	95	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 15.

All (95) 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)
3:A:422:PFZ:H92	4:A:715:HOH:O	1.47	1.11
2:A:417:HEM:ND	3:A:422:PFZ:H18	1.67	1.10
1:A:295:VAL:HG13	3:A:422:PFZ:H13	1.37	1.05
3:A:422:PFZ:CL21	3:A:422:PFZ:H41	1.96	1.02
2:A:417:HEM:C1D	3:A:422:PFZ:H18	1.97	0.98
3:A:422:PFZ:H62	4:A:712:HOH:O	1.64	0.96
1:A:87:PHE:CZ	3:A:422:PFZ:H103	2.04	0.92
1:A:252:THR:HG21	3:A:422:PFZ:H20	1.57	0.86
3:A:422:PFZ:CL21	3:A:422:PFZ:C4	2.60	0.86
1:A:87:PHE:CE1	3:A:422:PFZ:H103	2.20	0.77
1:A:87:PHE:CZ	3:A:422:PFZ:H71	2.27	0.70
1:A:252:THR:HG21	3:A:422:PFZ:C20	2.22	0.69
1:A:297:ASP:CG	3:A:422:PFZ:CL22	2.73	0.67
1:A:191[B]:MET:HE2	1:A:196:ALA:HA	1.75	0.66
1:A:98:PHE:HD2	3:A:422:PFZ:H82	1.60	0.66
1:A:181:THR:HG23	1:A:247:VAL:HG13	1.79	0.65
1:A:295:VAL:HG13	3:A:422:PFZ:C13	2.21	0.64
2:A:417:HEM:C4D	3:A:422:PFZ:H18	2.34	0.62
1:A:231:ARG:HB2	1:A:232:PRO:HD2	1.82	0.61
1:A:121:MET:HB2	1:A:122:PRO:HD3	1.81	0.61
1:A:158:PHE:HB3	1:A:159:PRO:HD3	1.84	0.60
1:A:99:ILE:HG12	1:A:103:MET:HE3	1.83	0.59
1:A:325:SER:O	1:A:331:GLU:HG3	2.04	0.57
1:A:251:ASP:O	1:A:255:ASN:ND2	2.38	0.57
1:A:149:ASN:ND2	1:A:402:LEU:H	2.03	0.56
1:A:128:GLU:HG3	1:A:365[A]:ARG:HH11	1.70	0.56
1:A:108:GLN:HE22	1:A:354:SER:HB2	1.69	0.56
1:A:295:VAL:HG11	2:A:417:HEM:HMA2	1.88	0.56
1:A:214:LYS:O	1:A:214:LYS:HD3	2.06	0.55
1:A:170:PRO:HG2	1:A:173:ASP:OD2	2.08	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:276:GLU:C	1:A:278:PRO:HD3	2.33	0.54
1:A:101:THR:OG1	3:A:422:PFZ:H81	2.09	0.53
1:A:150:PHE:CZ	1:A:261:MET:HG3	2.44	0.53
1:A:87:PHE:CE2	3:A:422:PFZ:H103	2.43	0.52
1:A:30:ASN:ND2	4:A:555:HOH:O	2.41	0.52
1:A:184:MET:HE3	4:A:713:HOH:O	2.09	0.52
1:A:205:ILE:HB	1:A:206:PRO:HD3	1.92	0.52
1:A:208:ILE:O	1:A:212:ARG:HG3	2.10	0.51
1:A:87:PHE:CD2	1:A:395[1]:ILE:HD13	2.46	0.51
1:A:114:LEU:O	1:A:117:GLN:HB2	2.10	0.51
1:A:56:THR:O	1:A:61:GLY:HA2	2.10	0.51
1:A:14:LEU:HD11	1:A:18:VAL:HG11	1.93	0.50
1:A:295:VAL:CG1	3:A:422:PFZ:H13	2.26	0.50
1:A:350:PHE:HB3	1:A:357:CYS:HB3	1.93	0.50
1:A:219:ALA:O	1:A:223:VAL:HG23	2.12	0.50
1:A:134:LEU:O	1:A:138:LEU:HB2	2.13	0.49
1:A:19:PRO:HG2	1:A:22:LEU:HD12	1.93	0.49
3:A:422:PFZ:C4	3:A:422:PFZ:H17	2.43	0.49
1:A:319:LEU:C	1:A:321:PRO:HD3	2.39	0.48
1:A:150:PHE:CE2	1:A:155:ALA:HB2	2.49	0.48
1:A:149:ASN:HD21	1:A:402:LEU:H	1.61	0.47
1:A:277:ARG:N	1:A:278:PRO:HD3	2.29	0.47
1:A:297:ASP:OD1	3:A:422:PFZ:CL22	2.69	0.47
1:A:183:GLN:OE1	1:A:188:ASP:HB3	2.15	0.47
1:A:181:THR:HG22	1:A:251:ASP:HB2	1.96	0.47
1:A:122:PRO:HD2	4:A:701:HOH:O	2.14	0.47
1:A:233:ILE:HG12	1:A:237:GLU:HB2	1.97	0.46
1:A:121:MET:CB	1:A:122:PRO:HD3	2.45	0.46
1:A:14:LEU:HD11	1:A:18:VAL:CG1	2.45	0.46
1:A:78:TYR:CE2	1:A:105:PRO:HG2	2.51	0.46
1:A:244:LEU:HD12	3:A:422:PFZ:H61	1.97	0.46
1:A:317:GLN:NE2	4:A:647:HOH:O	2.49	0.46
1:A:302:THR:O	1:A:314:LYS:HE3	2.15	0.45
1:A:108:GLN:HG3	1:A:355:HIS:CE1	2.52	0.45
1:A:114:LEU:HD23	1:A:241:MET:HE1	1.99	0.45
3:A:422:PFZ:C4	3:A:422:PFZ:C17	2.95	0.45
1:A:134:LEU:HD23	1:A:162:ILE:HG13	1.98	0.45
1:A:127:LEU:HD11	1:A:166:LEU:HD13	1.99	0.44
1:A:17:HIS:CD2	1:A:313:LYS:HG3	2.53	0.44
1:A:242:CYS:O	1:A:246:LEU:HD23	2.18	0.44
1:A:295:VAL:CG1	3:A:422:PFZ:CL22	3.03	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:181:THR:CG2	1:A:251:ASP:HB2	2.48	0.43
1:A:319:LEU:O	1:A:321:PRO:HD3	2.18	0.43
1:A:362:LEU:O	1:A:366:GLU:HG3	2.18	0.43
1:A:236:ASP:O	1:A:240:ARG:HG3	2.18	0.43
1:A:203:TYR:O	1:A:206:PRO:HD2	2.18	0.43
1:A:395[2]:ILE:HD13	4:A:714:HOH:O	2.19	0.43
3:A:422:PFZ:C9	4:A:715:HOH:O	2.31	0.43
1:A:49:ASN:HD22	1:A:49:ASN:H	1.67	0.43
1:A:40:GLU:HG3	1:A:336:MET:CE	2.48	0.42
1:A:294:LEU:HD23	1:A:294:LEU:H	1.85	0.42
1:A:83:SER:HB3	1:A:101:THR:O	2.20	0.42
1:A:191[B]:MET:HE2	1:A:196:ALA:CA	2.47	0.42
1:A:63:TRP:HB2	1:A:318:ILE:HG13	2.02	0.42
1:A:101:THR:HG21	3:A:422:PFZ:CL21	2.56	0.42
1:A:395[2]:ILE:CD1	4:A:714:HOH:O	2.68	0.41
1:A:211:ARG:NH1	1:A:218:ASP:OD2	2.54	0.41
1:A:407:ASP:HA	1:A:408:PRO:HD3	1.85	0.41
1:A:350:PHE:HB3	1:A:357:CYS:CB	2.51	0.41
1:A:132:GLN:NE2	1:A:373:GLU:OE2	2.53	0.41
1:A:114:LEU:HD23	1:A:241:MET:CE	2.51	0.41
1:A:200:LEU:O	1:A:203:TYR:HB3	2.21	0.40
1:A:40:GLU:HG3	1:A:336:MET:HE2	2.03	0.40
1:A:70:LEU:HD22	1:A:308:HIS:CD2	2.56	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	407/414 (98%)	390 (96%)	15 (4%)	2 (0%)	24 10

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	120	GLY
1	A	150	PHE

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	354/358 (99%)	343 (97%)	11 (3%)	35 13

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

Mol	Chain	Res	Type
1	A	49	ASN
1	A	79	ARG
1	A	86	PRO
1	A	96	TYR
1	A	108	GLN
1	A	184	MET
1	A	195	GLU
1	A	222	ILE
1	A	235	SER
1	A	272	GLN
1	A	276	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	21	HIS
1	A	30	ASN
1	A	39	GLN
1	A	46	GLN
1	A	49	ASN
1	A	108	GLN
1	A	110	GLN
1	A	116	ASN
1	A	117	GLN

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Mol	Chain	Res	Type
1	A	129	ASN
1	A	149	ASN
1	A	225	ASN
1	A	229	ASN
1	A	317	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](#)

2 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	PFZ	A	422	2	23,23,23	1.18	2 (8%)	28,31,31	1.27	3 (10%)
2	HEM	A	417	1,3	50,50,50	1.42	7 (14%)	67,82,82	1.27	7 (10%)

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	PFZ	A	422	2	-	16/18/18/18	0/2/2/2
2	HEM	A	417	1,3	-	2/14/54/54	-

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	417	HEM	FE-NB	3.26	2.04	1.94
2	A	417	HEM	FE-NC	3.09	2.05	1.95
2	A	417	HEM	FE-NA	3.04	2.05	1.95
3	A	422	PFZ	C5-N11	3.02	1.51	1.46
2	A	417	HEM	C4A-C3A	2.56	1.49	1.43
2	A	417	HEM	CAC-C3C	2.51	1.54	1.47
2	A	417	HEM	C2A-C3A	-2.24	1.33	1.38
3	A	422	PFZ	C20-N19	2.20	1.37	1.32
2	A	417	HEM	CMB-C2B	2.09	1.55	1.50

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	417	HEM	O1D-CGD-CBD	-3.24	112.81	123.09
2	A	417	HEM	CMA-C3A-C4A	-3.05	120.78	125.42
2	A	417	HEM	CMA-C3A-C2A	3.02	132.03	125.62
2	A	417	HEM	C3B-C2B-C1B	2.81	108.52	106.41
2	A	417	HEM	CBD-CAD-C3D	2.34	119.00	112.53
3	A	422	PFZ	C18-N19-C20	2.21	109.70	105.47
2	A	417	HEM	CBC-CAC-C3C	-2.17	116.66	127.53
3	A	422	PFZ	C5-N11-C20	2.17	129.74	125.75
2	A	417	HEM	C3D-C4D-ND	-2.12	107.84	110.17
3	A	422	PFZ	C6-C4-C1	-2.02	113.42	116.01

There are no chirality outliers.

All (18) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	417	HEM	C2C-C3C-CAC-CBC
2	A	417	HEM	C4C-C3C-CAC-CBC
3	A	422	PFZ	O2-C1-C3-C12
3	A	422	PFZ	O2-C1-C3-C16
3	A	422	PFZ	C4-C1-C3-C12
3	A	422	PFZ	C4-C1-C3-C16
3	A	422	PFZ	C5-C1-C3-C12
3	A	422	PFZ	C5-C1-C3-C16

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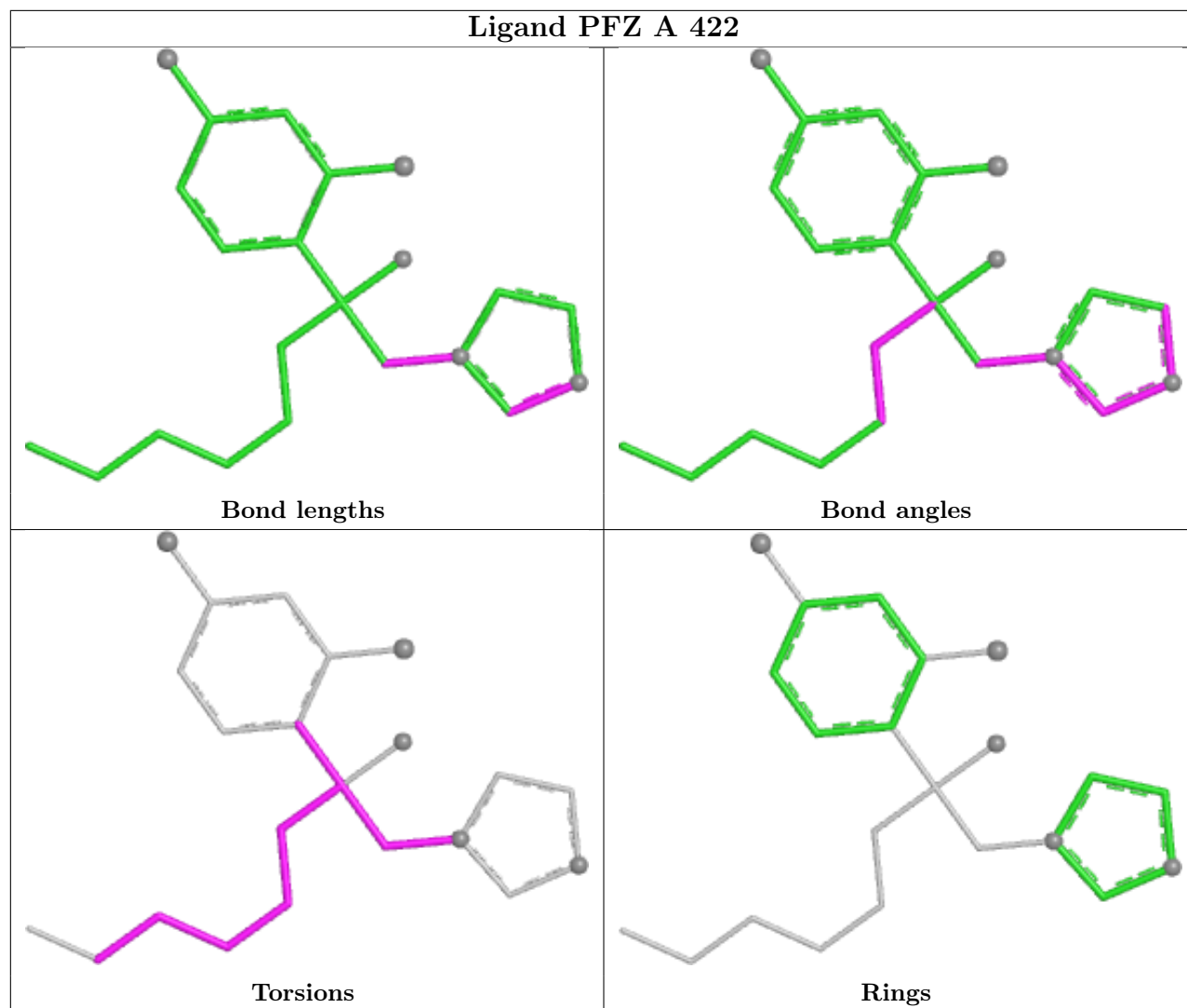
Mol	Chain	Res	Type	Atoms
3	A	422	PFZ	O2-C1-C4-C6
3	A	422	PFZ	C3-C1-C4-C6
3	A	422	PFZ	C5-C1-C4-C6
3	A	422	PFZ	C1-C5-N11-C20
3	A	422	PFZ	O2-C1-C5-N11
3	A	422	PFZ	C4-C6-C7-C8
3	A	422	PFZ	C7-C8-C9-C10
3	A	422	PFZ	C6-C7-C8-C9
3	A	422	PFZ	C1-C5-N11-C17
3	A	422	PFZ	C1-C4-C6-C7

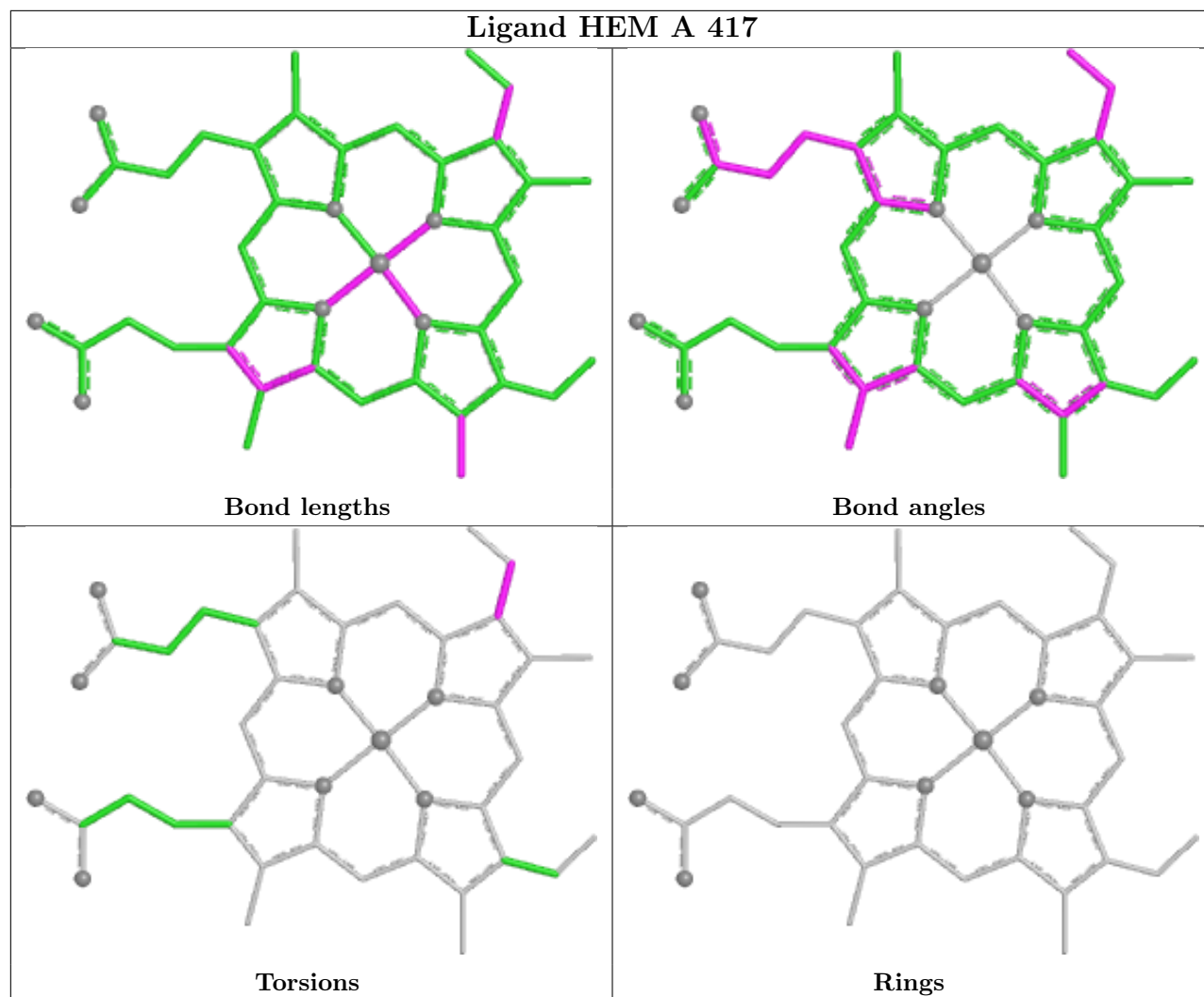
There are no ring outliers.

2 monomers are involved in 27 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	422	PFZ	26	0
2	A	417	HEM	4	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

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.