



# Full wwPDB X-ray Structure Validation Report ⓘ

May 22, 2020 – 11:52 pm BST

PDB ID : 2WX2  
Title : X-RAY STRUCTURE OF CYP51 FROM THE HUMAN PATHOGEN TRY-  
PANOSOMA CRUZI IN COMPLEX WITH FLUCONAZOLE  
Authors : Chen, C.-K.; Leung, S.S.F.; Guilbert, C.; Jacobson, M.; McKerrow, J.H.;  
Podust, L.M.  
Deposited on : 2009-10-31  
Resolution : 2.27 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](mailto: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.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.11  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.11

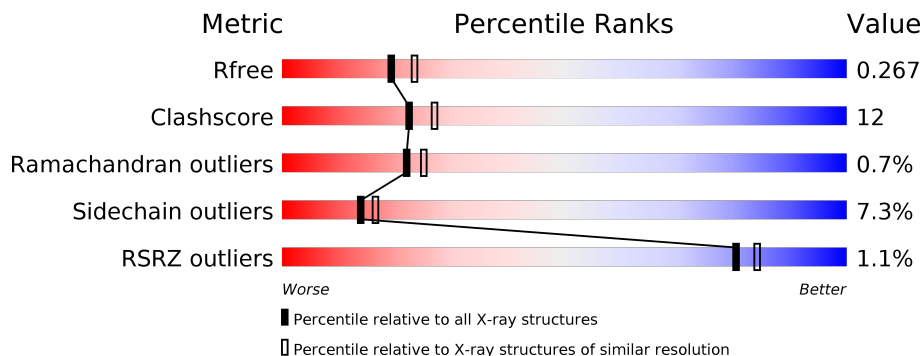
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.27 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))
$R_{free}$	130704	6980 (2.30-2.26)
Clashscore	141614	7711 (2.30-2.26)
Ramachandran outliers	138981	7597 (2.30-2.26)
Sidechain outliers	138945	7598 (2.30-2.26)
RSRZ outliers	127900	6849 (2.30-2.26)

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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	473	
1	B	473	

## 2 Entry composition i

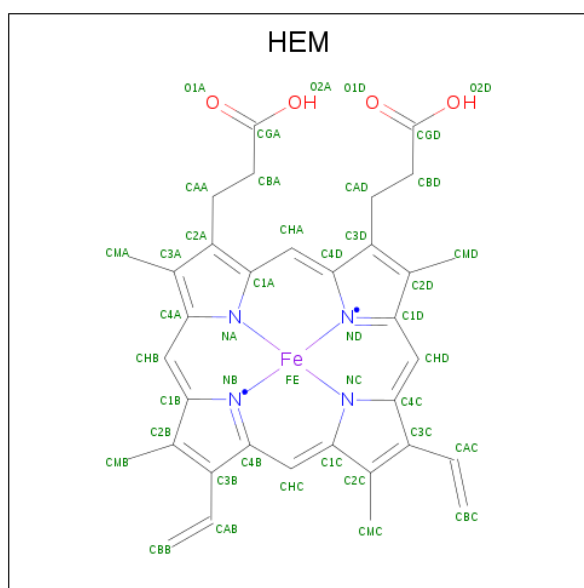
There are 4 unique types of molecules in this entry. The entry contains 7513 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 LANOSTEROL 14-ALPHA-DEMETHYLASE.

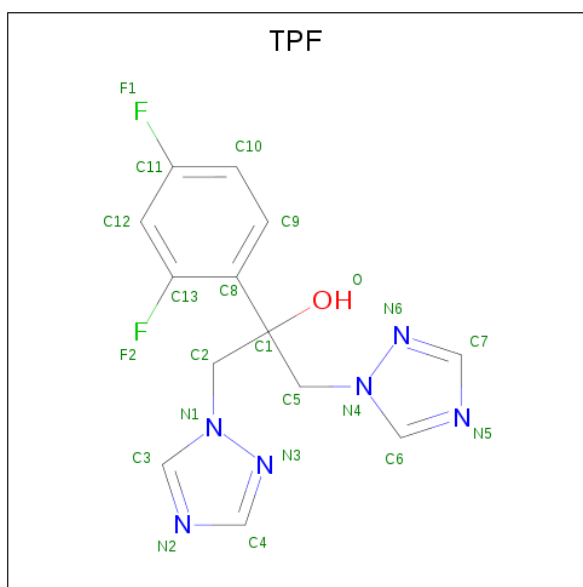
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	441	Total	C	N	O	S	0	1	0
			3469	2226	604	612	27			
1	B	444	Total	C	N	O	S	0	0	0
			3485	2239	599	620	27			

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



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

- Molecule 3 is 2-(2,4-DIFLUOROPHENYL)-1,3-DI(1H-1,2,4-TRIAZOL-1-YL)PROPAN-2-OL (three-letter code: TPF) (formula:  $C_{13}H_{12}F_2N_6O$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	F	N	O		
3	A	1	22	13	2	6	1	0	0
3	B	1	22	13	2	6	1	0	0

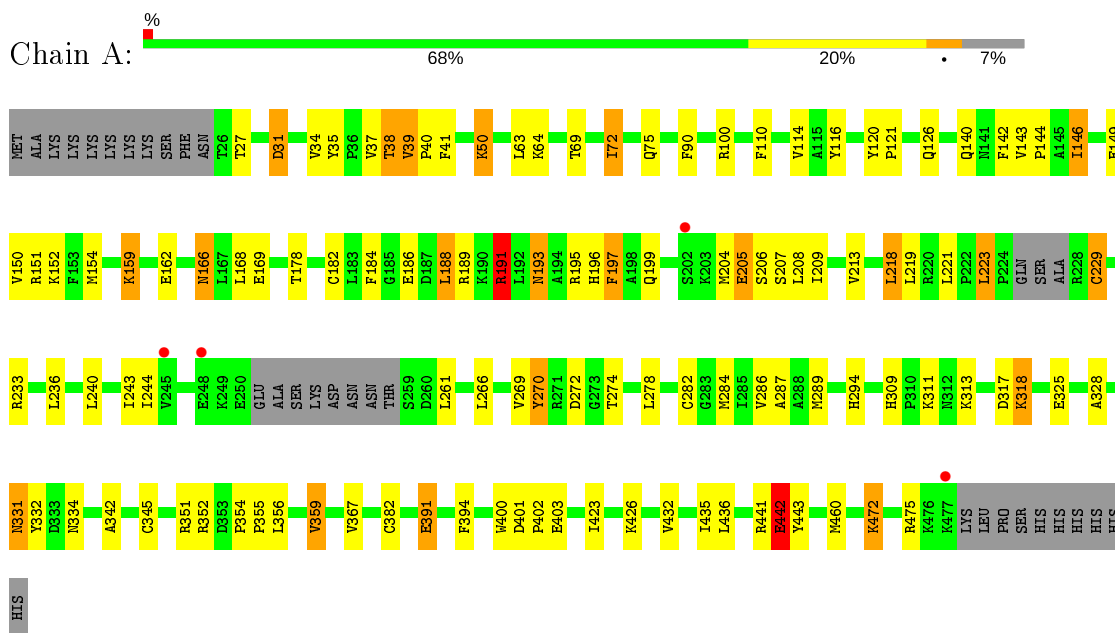
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	A	217	217	217	0	0
4	B	212	212	212	0	0

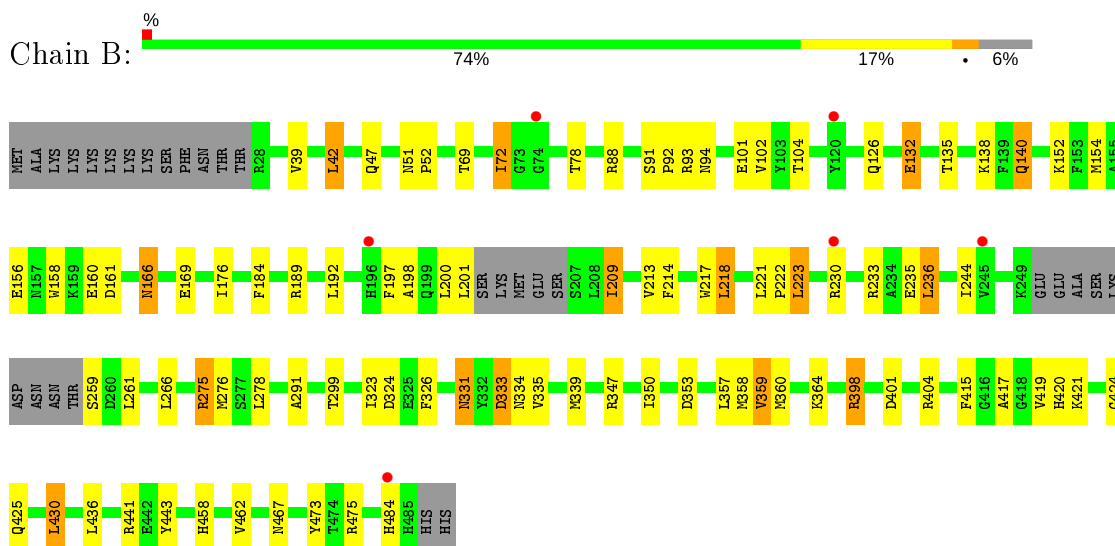
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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: LANOSTEROL 14-ALPHA-DEMETHYLASE



- Molecule 1: LANOSTEROL 14-ALPHA-DEMETHYLASE



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	69.99Å 101.47Å 74.73Å 90.00° 111.63° 90.00°	Depositor
Resolution (Å)	69.50 – 2.27 24.25 – 2.27	Depositor EDS
% Data completeness (in resolution range)	89.1 (69.50-2.27) 89.2 (24.25-2.27)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.01 (at 2.26Å)	Xtrriage
Refinement program	REFMAC 5.5.0072	Depositor
R, $R_{free}$	0.193 , 0.273 0.189 , 0.267	Depositor DCC
$R_{free}$ test set	1944 reflections (4.86%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	25.0	Xtrriage
Anisotropy	0.093	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 36.7	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	7513	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	23.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.75% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: TPF, 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	0.77	1/3553 (0.0%)	0.85	6/4816 (0.1%)
1	B	0.75	0/3571	0.80	3/4844 (0.1%)
All	All	0.76	1/7124 (0.0%)	0.83	9/9660 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	442	GLU	CG-CD	-5.76	1.43	1.51

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	223	LEU	CA-CB-CG	6.46	130.16	115.30
1	A	191	ARG	NE-CZ-NH1	5.86	123.23	120.30
1	B	333	ASP	CB-CG-OD1	5.72	123.45	118.30
1	A	223	LEU	CA-CB-CG	5.56	128.09	115.30
1	A	359	VAL	CB-CA-C	5.51	121.87	111.40
1	A	352	ARG	NE-CZ-NH2	-5.33	117.64	120.30
1	B	436	LEU	CA-CB-CG	5.24	127.35	115.30
1	A	31	ASP	N-CA-CB	5.13	119.83	110.60
1	A	401	ASP	CB-CG-OD1	5.08	122.87	118.30

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	3469	0	3419	86	0
1	B	3485	0	3426	74	0
2	A	43	0	30	7	0
2	B	43	0	30	3	0
3	A	22	0	12	1	0
3	B	22	0	12	1	0
4	A	217	0	0	4	0
4	B	212	0	0	7	0
All	All	7513	0	6929	162	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 12.

All (162) 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:191:ARG:HG2	1:A:191:ARG:HH11	1.21	1.04
1:B:184:PHE:O	1:B:189:ARG:NH1	1.93	1.02
1:B:259:SER:HA	4:B:2085:HOH:O	1.59	1.00
1:B:331:ASN:H	1:B:334:ASN:HD22	1.16	0.89
1:A:331:ASN:H	1:A:334:ASN:HD22	1.17	0.86
1:A:193:ASN:ND2	1:A:196:HIS:H	1.74	0.85
1:A:269:VAL:O	1:A:274:THR:O	1.93	0.85
1:B:331:ASN:H	1:B:334:ASN:ND2	1.74	0.85
1:A:114:VAL:O	1:A:126:GLN:OE1	1.93	0.84
1:B:192:LEU:HD21	1:B:197:PHE:HD1	1.41	0.83
1:A:309:HIS:CE1	1:A:311:LYS:HG2	2.14	0.82
1:A:191:ARG:HG2	1:A:191:ARG:NH1	1.92	0.82
2:A:1450:HEM:HBC2	2:A:1450:HEM:HMC2	1.63	0.80
1:B:39:VAL:HA	1:B:72:ILE:HD11	1.60	0.80
1:A:309:HIS:ND1	1:A:311:LYS:HG2	1.98	0.79
1:A:331:ASN:ND2	1:A:334:ASN:H	1.85	0.73
1:A:282:CYS:O	1:A:286:VAL:HG23	1.88	0.73
1:B:154:MET:HE2	1:B:443:TYR:HE2	1.54	0.72
1:B:154:MET:CE	1:B:443:TYR:HE2	2.03	0.72
1:B:192:LEU:HD21	1:B:197:PHE:CD1	2.23	0.71
1:A:184:PHE:O	1:A:189:ARG:NH1	2.24	0.70
1:A:38:THR:HG22	1:A:41:PHE:H	1.58	0.69
1:B:443:TYR:OH	1:B:475:ARG:NH2	2.26	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:149:GLU:HB2	1:A:178:THR:HG22	1.76	0.67
1:A:331:ASN:H	1:A:334:ASN:ND2	1.91	0.66
1:B:347:ARG:HD3	4:B:2134:HOH:O	1.96	0.66
1:B:47:GLN:O	1:B:458:HIS:HE1	1.79	0.66
1:A:142:PHE:O	1:A:146:ILE:HG23	1.95	0.66
1:A:38:THR:CG2	1:A:41:PHE:H	2.10	0.65
1:B:217:TRP:CZ3	1:B:218:LEU:HD13	2.32	0.65
1:B:401:ASP:O	1:B:404:ARG:HB3	1.97	0.65
1:A:146:ILE:HG22	1:A:182:CYS:SG	2.36	0.64
1:A:244:ILE:HD12	1:A:278:LEU:HD13	1.80	0.64
1:A:233[B]:ARG:HB3	1:A:233[B]:ARG:HH11	1.62	0.64
1:A:191:ARG:CG	1:A:191:ARG:HH11	2.04	0.64
1:A:240:LEU:O	1:A:244:ILE:HG13	1.98	0.63
1:A:166:ASN:HD22	1:A:169:GLU:H	1.45	0.62
1:A:309:HIS:ND1	1:A:311:LYS:CG	2.62	0.62
1:A:269:VAL:O	1:A:270:TYR:HB2	1.97	0.62
1:A:272:ASP:HA	1:B:140:GLN:NE2	2.14	0.62
1:A:197:PHE:CZ	1:A:289:MET:HG3	2.34	0.62
1:B:291:ALA:O	3:B:1460:TPF:HC7	2.00	0.61
1:A:391:GLU:HG3	4:A:2164:HOH:O	2.00	0.60
1:B:326:PHE:HE2	1:B:339:MET:HE2	1.67	0.60
1:B:42:LEU:HD13	1:B:214:PHE:CZ	2.37	0.60
1:A:294:HIS:HB2	4:A:2115:HOH:O	2.02	0.59
1:A:331:ASN:HD22	1:A:334:ASN:H	1.50	0.59
1:A:318:LYS:HE2	1:A:403:GLU:OE2	2.03	0.58
1:B:324:ASP:OD2	4:B:2119:HOH:O	2.16	0.58
1:B:152:LYS:HE2	1:B:156:GLU:OE2	2.03	0.58
1:A:269:VAL:O	1:A:270:TYR:CB	2.52	0.58
1:A:342:ALA:O	1:A:345:CYS:HB2	2.04	0.57
2:A:1450:HEM:HBC2	2:A:1450:HEM:CMC	2.33	0.56
1:A:193:ASN:HD21	1:A:196:HIS:H	1.52	0.56
1:A:204:MET:C	1:A:206:SER:H	2.08	0.56
1:B:353:ASP:OD1	1:B:398:ARG:NH2	2.39	0.55
1:A:432:VAL:O	1:A:436:LEU:HG	2.06	0.55
1:B:197:PHE:HE2	1:B:201:LEU:HD21	1.71	0.55
3:A:1460:TPF:F2	3:A:1460:TPF:HC52	1.97	0.55
1:B:331:ASN:HD21	1:B:333:ASP:HB2	1.72	0.55
1:A:233[B]:ARG:NH1	1:A:233[B]:ARG:HB3	2.20	0.55
1:B:230:ARG:HG3	1:B:233:ARG:NH2	2.22	0.55
1:A:150:VAL:O	1:A:154:MET:HG3	2.07	0.55
1:B:244:ILE:HG12	1:B:266:LEU:HD11	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:166:ASN:HD22	1:B:169:GLU:H	1.55	0.54
1:B:154:MET:CE	1:B:443:TYR:CE2	2.89	0.54
1:A:313:LYS:HD3	1:A:317:ASP:OD1	2.08	0.54
1:A:34:VAL:HG22	1:A:69:THR:HB	1.89	0.53
1:B:331:ASN:ND2	1:B:333:ASP:H	2.06	0.53
1:A:75:GLN:OE1	1:A:218:LEU:HD22	2.08	0.53
1:A:218:LEU:C	1:A:219:LEU:O	2.44	0.53
1:B:424:GLY:HA3	2:B:1450:HEM:C3C	2.44	0.53
1:A:193:ASN:HD22	1:A:196:HIS:H	1.51	0.53
1:A:442:GLU:HB2	1:A:443:TYR:CD2	2.44	0.53
1:A:356:LEU:HD12	2:A:1450:HEM:C3A	2.44	0.52
1:B:417:ALA:HA	1:B:421:LYS:HG3	1.91	0.52
1:A:331:ASN:N	1:A:334:ASN:HD22	1.97	0.52
1:B:221:LEU:HB3	1:B:222:PRO:CD	2.40	0.52
1:A:90:PHE:HE2	1:A:382:CYS:HG	1.59	0.51
1:A:209:ILE:O	1:A:213:VAL:HG22	2.12	0.50
1:B:335:VAL:HG12	1:B:430:LEU:HD12	1.93	0.50
1:A:331:ASN:C	1:A:331:ASN:HD22	2.14	0.50
1:A:244:ILE:HD12	1:A:278:LEU:CD1	2.41	0.50
1:B:326:PHE:CE2	1:B:339:MET:CE	2.95	0.50
1:A:114:VAL:O	1:A:284:MET:HE3	2.12	0.50
1:B:331:ASN:N	1:B:334:ASN:HD22	1.98	0.50
1:A:35:TYR:HB2	1:A:63:LEU:HD11	1.94	0.49
1:B:135:THR:HA	1:B:138:LYS:HD2	1.94	0.49
1:B:326:PHE:CE2	1:B:339:MET:HE2	2.46	0.49
1:A:39:VAL:HG22	1:A:40:PRO:HD3	1.93	0.49
1:B:158:TRP:HB3	1:B:473:TYR:CZ	2.47	0.49
1:A:143:VAL:HB	1:A:144:PRO:HD3	1.95	0.48
1:A:38:THR:HG22	1:A:41:PHE:HB3	1.94	0.48
1:B:424:GLY:HA3	2:B:1450:HEM:C2C	2.47	0.48
1:B:420:HIS:HD1	2:B:1450:HEM:CGD	2.24	0.48
1:A:151:ARG:NH2	1:A:328:ALA:O	2.46	0.48
1:A:150:VAL:HG21	1:A:435:ILE:HG12	1.96	0.48
1:B:358:MET:HG3	1:B:359:VAL:N	2.28	0.47
1:A:37:VAL:O	1:A:72:ILE:O	2.32	0.47
1:B:200:LEU:HD12	1:B:200:LEU:N	2.29	0.47
1:B:154:MET:HE2	1:B:443:TYR:CE2	2.41	0.47
1:B:404:ARG:HG3	4:B:2162:HOH:O	2.15	0.47
1:A:204:MET:C	1:A:206:SER:N	2.68	0.47
1:B:200:LEU:HD23	1:B:236:LEU:HD12	1.96	0.47
1:B:275:ARG:HD3	4:B:2093:HOH:O	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:75:GLN:CD	1:A:218:LEU:HD22	2.36	0.46
1:A:100:ARG:HG3	1:A:116:TYR:O	2.15	0.46
1:B:200:LEU:CD1	1:B:200:LEU:N	2.78	0.46
1:B:415:PHE:CD1	1:B:425:GLN:HG3	2.49	0.46
1:B:197:PHE:CE2	1:B:201:LEU:HD21	2.50	0.46
1:B:102:VAL:HB	1:B:360:MET:HE3	1.99	0.45
1:B:154:MET:HE3	1:B:443:TYR:CE2	2.52	0.45
1:B:69:THR:HA	1:B:78:THR:HA	1.98	0.45
1:B:88:ARG:HA	1:B:88:ARG:HD3	1.68	0.45
1:B:94:ASN:ND2	1:B:94:ASN:H	2.15	0.45
1:A:204:MET:HA	1:A:229:CYS:HB2	1.98	0.45
1:A:356:LEU:HD12	2:A:1450:HEM:CMA	2.47	0.45
1:A:38:THR:HG22	1:A:41:PHE:CB	2.47	0.45
1:A:50:LYS:HD2	1:A:50:LYS:HA	1.73	0.45
1:A:184:PHE:CE1	1:A:261:LEU:HD23	2.51	0.45
1:B:323:ILE:HD12	1:B:441:ARG:HG3	1.99	0.45
1:B:101:GLU:OE1	1:B:364:LYS:NZ	2.45	0.45
1:B:326:PHE:HE2	1:B:339:MET:CE	2.31	0.44
1:B:244:ILE:HD12	1:B:278:LEU:HD13	1.99	0.44
1:A:159:LYS:HB2	1:A:159:LYS:NZ	2.29	0.44
1:A:193:ASN:C	1:A:193:ASN:HD22	2.21	0.44
1:A:191:ARG:HD3	1:A:243:ILE:HG12	2.00	0.43
1:B:154:MET:HG2	1:B:158:TRP:CE3	2.52	0.43
1:A:27:THR:HG22	4:A:2004:HOH:O	2.17	0.43
1:A:354:PRO:HA	1:A:355:PRO:HD3	1.84	0.43
1:B:126:GLN:HG2	1:B:276:MET:CE	2.48	0.43
1:B:91:SER:HB3	1:B:92:PRO:HD3	1.99	0.43
1:B:51:ASN:HA	1:B:52:PRO:HD3	1.80	0.43
1:A:146:ILE:HD11	1:A:435:ILE:HG13	2.01	0.43
1:B:209:ILE:O	1:B:213:VAL:HB	2.19	0.42
1:A:205:GLU:HA	1:A:208:LEU:HB2	2.01	0.42
1:B:419:VAL:HG23	1:B:420:HIS:CD2	2.54	0.42
1:B:218:LEU:HA	1:B:218:LEU:HD12	1.75	0.42
1:A:367:VAL:HG22	4:A:2040:HOH:O	2.18	0.42
1:A:332:TYR:CE2	1:B:93:ARG:HD2	2.55	0.42
2:A:1450:HEM:CMB	2:A:1450:HEM:HBB2	2.50	0.42
1:B:132:GLU:O	1:B:135:THR:HG22	2.20	0.42
1:B:192:LEU:CD2	1:B:197:PHE:HD1	2.23	0.42
1:A:188:LEU:HD22	1:A:243:ILE:HD13	2.02	0.41
1:A:120:TYR:CE1	1:A:121:PRO:HB3	2.55	0.41
1:A:110:PHE:CE1	1:A:287:ALA:HA	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:318:LYS:CE	1:A:403:GLU:OE2	2.68	0.41
1:B:176:ILE:HG12	1:B:198:ALA:HB2	2.01	0.41
1:B:299:THR:HG21	1:B:350:ILE:HD11	2.02	0.41
1:B:347:ARG:NH2	4:B:2135:HOH:O	2.53	0.41
1:A:356:LEU:HD12	2:A:1450:HEM:HMA2	2.02	0.41
1:A:162:GLU:OE2	1:A:472:LYS:HE2	2.20	0.41
1:A:38:THR:HG23	1:A:40:PRO:HD2	2.03	0.41
1:A:159:LYS:CB	1:A:159:LYS:NZ	2.84	0.41
1:A:351:ARG:HD3	1:A:394:PHE:CD1	2.56	0.41
1:B:161:ASP:HB3	4:B:2208:HOH:O	2.20	0.41
1:B:214:PHE:O	1:B:217:TRP:HB3	2.20	0.41
1:B:357:LEU:HD11	1:B:462:VAL:HG21	2.02	0.40
1:A:400:TRP:CH2	1:A:402:PRO:HG3	2.56	0.40
1:B:244:ILE:CD1	1:B:278:LEU:HD13	2.51	0.40
1:A:72:ILE:HG21	1:A:218:LEU:HD11	2.02	0.40
1:A:423:ILE:HD11	2:A:1450:HEM:HMD2	2.03	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	436/473 (92%)	415 (95%)	17 (4%)	4 (1%)	17	18
1	B	438/473 (93%)	415 (95%)	21 (5%)	2 (0%)	29	34
All	All	874/946 (92%)	830 (95%)	38 (4%)	6 (1%)	22	25

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	140	GLN
1	B	484	HIS

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Mol	Chain	Res	Type
1	A	205	GLU
1	A	270	TYR
1	B	140	GLN
1	A	72	ILE

### 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	364/415 (88%)	329 (90%)	35 (10%)	8	9
1	B	367/415 (88%)	349 (95%)	18 (5%)	25	33
All	All	731/830 (88%)	678 (93%)	53 (7%)	14	16

All (53) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	31	ASP
1	A	38	THR
1	A	39	VAL
1	A	50	LYS
1	A	64	LYS
1	A	146	ILE
1	A	152	LYS
1	A	159	LYS
1	A	166	ASN
1	A	168	LEU
1	A	186	GLU
1	A	188	LEU
1	A	191	ARG
1	A	193	ASN
1	A	195	ARG
1	A	197	PHE
1	A	199	GLN
1	A	207	SER
1	A	218	LEU

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
1	A	221	LEU
1	A	223	LEU
1	A	229	CYS
1	A	236	LEU
1	A	266	LEU
1	A	318	LYS
1	A	325	GLU
1	A	331	ASN
1	A	359	VAL
1	A	391	GLU
1	A	426	LYS
1	A	441	ARG
1	A	442	GLU
1	A	460	MET
1	A	472	LYS
1	A	475	ARG
1	B	42	LEU
1	B	72	ILE
1	B	104	THR
1	B	132	GLU
1	B	160	GLU
1	B	166	ASN
1	B	209	ILE
1	B	218	LEU
1	B	223	LEU
1	B	235	GLU
1	B	236	LEU
1	B	261	LEU
1	B	275	ARG
1	B	331	ASN
1	B	359	VAL
1	B	398	ARG
1	B	430	LEU
1	B	467	ASN

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

<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
1	A	75	GLN
1	A	166	ASN
1	A	193	ASN
1	A	293	GLN

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Mol	Chain	Res	Type
1	A	331	ASN
1	A	334	ASN
1	A	458	HIS
1	B	140	GLN
1	B	166	ASN
1	B	193	ASN
1	B	196	HIS
1	B	199	GLN
1	B	293	GLN
1	B	331	ASN
1	B	334	ASN
1	B	458	HIS

### 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 carbohydrates 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	TPF	B	1460	2	18,24,24	2.05	2 (11%)	24,34,34	2.76	9 (37%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	HEM	A	1450	1,3	27,50,50	2.37	10 (37%)	17,82,82	1.74	4 (23%)
2	HEM	B	1450	1,3	27,50,50	2.18	8 (29%)	17,82,82	1.96	5 (29%)
3	TPF	A	1460	2	18,24,24	2.04	4 (22%)	24,34,34	3.04	11 (45%)

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	TPF	B	1460	2	-	2/16/16/16	0/3/3/3
2	HEM	A	1450	1,3	-	4/6/54/54	-
2	HEM	B	1450	1,3	-	2/6/54/54	-
3	TPF	A	1460	2	-	3/16/16/16	0/3/3/3

All (24) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1460	TPF	C8-C13	6.29	1.48	1.38
3	B	1460	TPF	C8-C13	6.18	1.48	1.38
2	A	1450	HEM	C3D-C2D	5.36	1.53	1.37
2	B	1450	HEM	C3C-C2C	-4.87	1.33	1.40
2	A	1450	HEM	C3C-C2C	-4.80	1.33	1.40
2	A	1450	HEM	C3B-C2B	-4.72	1.33	1.40
3	B	1460	TPF	N3-N1	4.43	1.41	1.35
2	B	1450	HEM	C3D-C2D	4.35	1.50	1.37
2	B	1450	HEM	C3B-C2B	-4.26	1.34	1.40
2	A	1450	HEM	C3B-CAB	4.14	1.56	1.47
2	B	1450	HEM	C3C-CAC	3.43	1.54	1.47
3	A	1460	TPF	N3-N1	3.30	1.40	1.35
2	B	1450	HEM	C3B-CAB	3.27	1.54	1.47
3	A	1460	TPF	N6-N4	3.14	1.39	1.35
2	A	1450	HEM	C3C-CAC	3.12	1.54	1.47
2	A	1450	HEM	CAA-C2A	2.83	1.56	1.52
2	B	1450	HEM	C1C-C2C	2.52	1.48	1.42
3	A	1460	TPF	C1-C8	-2.30	1.50	1.53
2	B	1450	HEM	CAA-C2A	2.30	1.55	1.52
2	A	1450	HEM	C1C-C2C	2.20	1.47	1.42
2	A	1450	HEM	CAD-C3D	2.09	1.55	1.52
2	B	1450	HEM	C1D-ND	2.08	1.40	1.36
2	A	1450	HEM	C4A-NA	2.07	1.40	1.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1450	HEM	CMD-C2D	2.01	1.55	1.51

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1460	TPF	C3-N2-C4	7.01	110.24	102.34
3	A	1460	TPF	C3-N2-C4	6.68	109.86	102.34
3	A	1460	TPF	C6-N5-C7	6.62	109.80	102.34
3	B	1460	TPF	N2-C3-N1	-5.64	105.43	112.24
3	A	1460	TPF	N2-C3-N1	-5.35	105.77	112.24
3	B	1460	TPF	C6-N5-C7	5.11	108.10	102.34
2	B	1450	HEM	CAD-CBD-CGD	-5.08	104.14	112.67
3	B	1460	TPF	N5-C6-N4	-4.78	106.46	112.24
3	A	1460	TPF	N5-C6-N4	-4.78	106.46	112.24
3	A	1460	TPF	C12-C13-C8	-4.17	119.47	124.00
3	A	1460	TPF	C1-C8-C13	-3.95	119.59	122.84
2	B	1450	HEM	CAD-C3D-C2D	-3.64	116.80	127.25
3	B	1460	TPF	C1-C8-C13	-3.40	120.04	122.84
2	A	1450	HEM	CBD-CAD-C3D	3.32	118.59	112.48
3	A	1460	TPF	C2-C1-C8	-3.31	102.63	110.87
3	A	1460	TPF	C13-C12-C11	3.18	119.96	116.62
2	B	1450	HEM	CBD-CAD-C3D	2.90	117.82	112.48
3	A	1460	TPF	C5-C1-C2	2.89	115.02	109.02
2	A	1450	HEM	C4C-C3C-C2C	2.87	108.91	106.90
3	B	1460	TPF	C13-C12-C11	2.81	119.57	116.62
2	A	1450	HEM	C1D-C2D-C3D	-2.62	105.18	107.00
3	B	1460	TPF	C9-C10-C11	2.46	120.91	118.36
2	B	1450	HEM	CMB-C2B-C3B	2.42	129.21	124.68
3	B	1460	TPF	F1-C11-C10	2.37	122.58	118.54
3	A	1460	TPF	C9-C10-C11	2.20	120.64	118.36
3	A	1460	TPF	C9-C8-C13	2.19	118.82	116.10
2	B	1450	HEM	CMA-C3A-C4A	-2.12	125.21	128.46
2	A	1450	HEM	CAD-CBD-CGD	-2.11	109.13	112.67
3	B	1460	TPF	C9-C8-C1	2.00	123.42	120.95

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	1460	TPF	C1-C5-N4-C6
2	A	1450	HEM	C1A-C2A-CAA-CBA
2	A	1450	HEM	C3A-C2A-CAA-CBA

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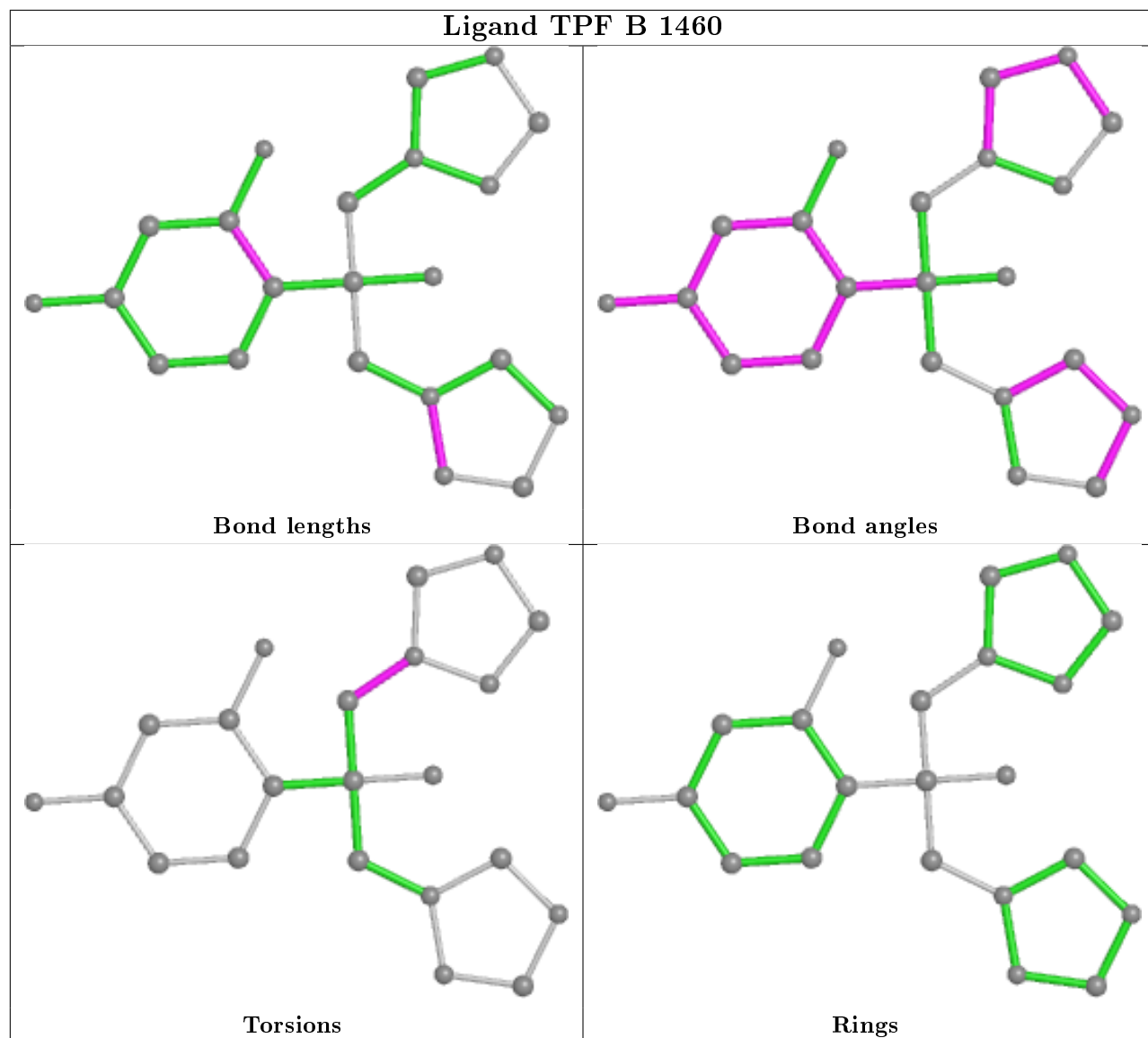
Mol	Chain	Res	Type	Atoms
2	A	1450	HEM	C2D-C3D-CAD-CBD
2	A	1450	HEM	C4D-C3D-CAD-CBD
2	B	1450	HEM	C2D-C3D-CAD-CBD
2	B	1450	HEM	C4D-C3D-CAD-CBD
3	A	1460	TPF	C1-C5-N4-C6
3	B	1460	TPF	C1-C5-N4-N6
3	A	1460	TPF	C1-C5-N4-N6
3	A	1460	TPF	C5-C1-C2-N1

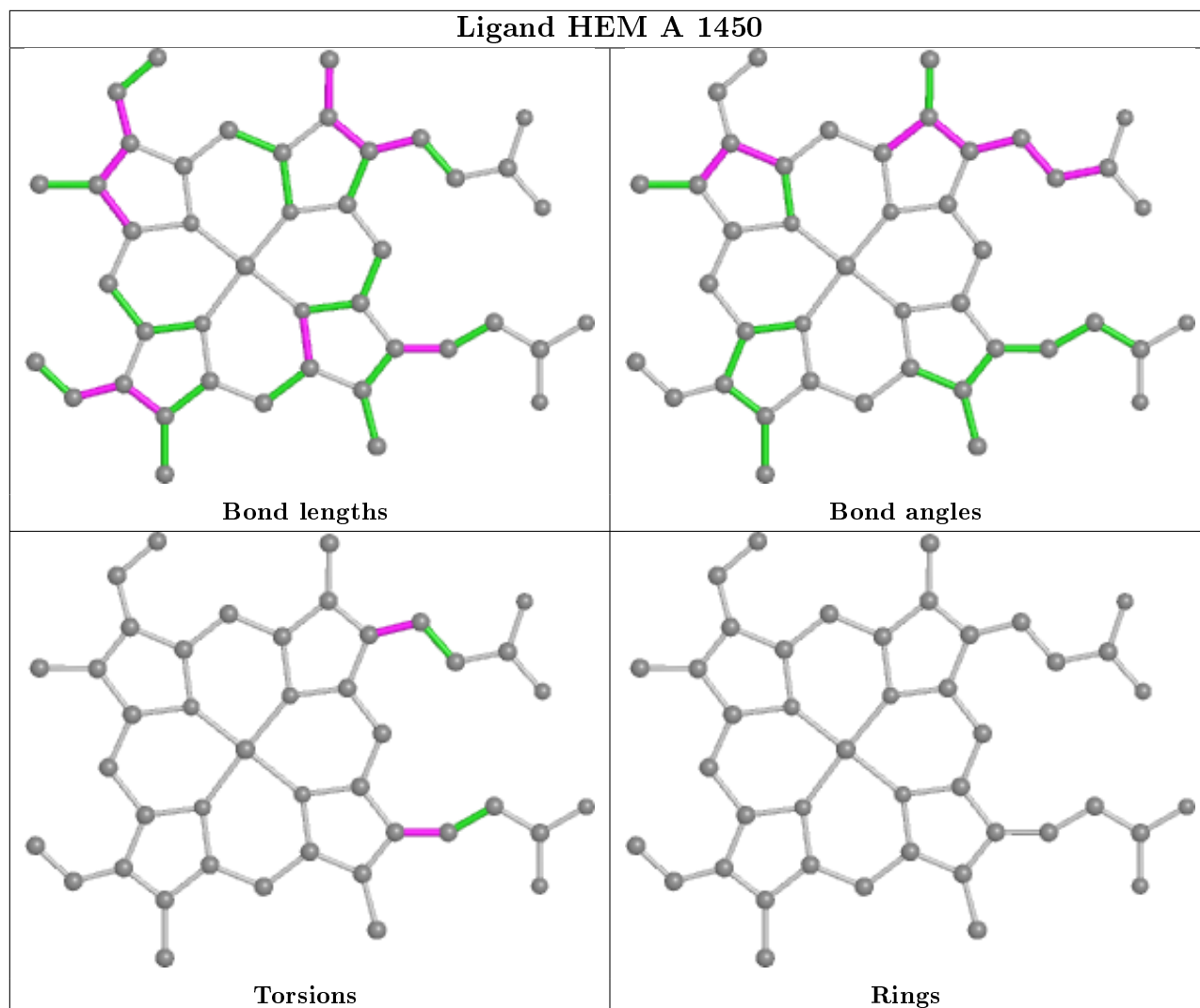
There are no ring outliers.

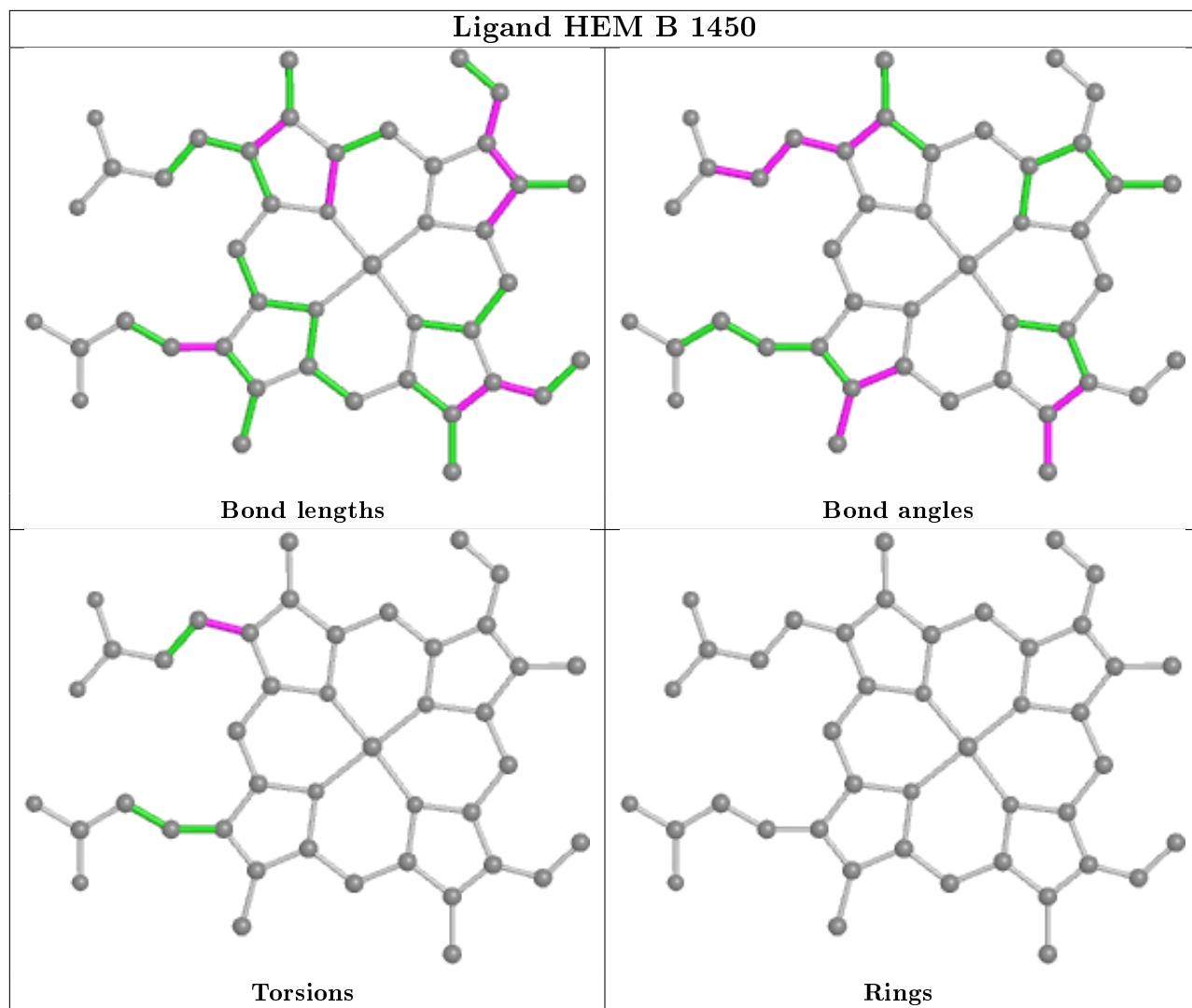
4 monomers are involved in 12 short contacts:

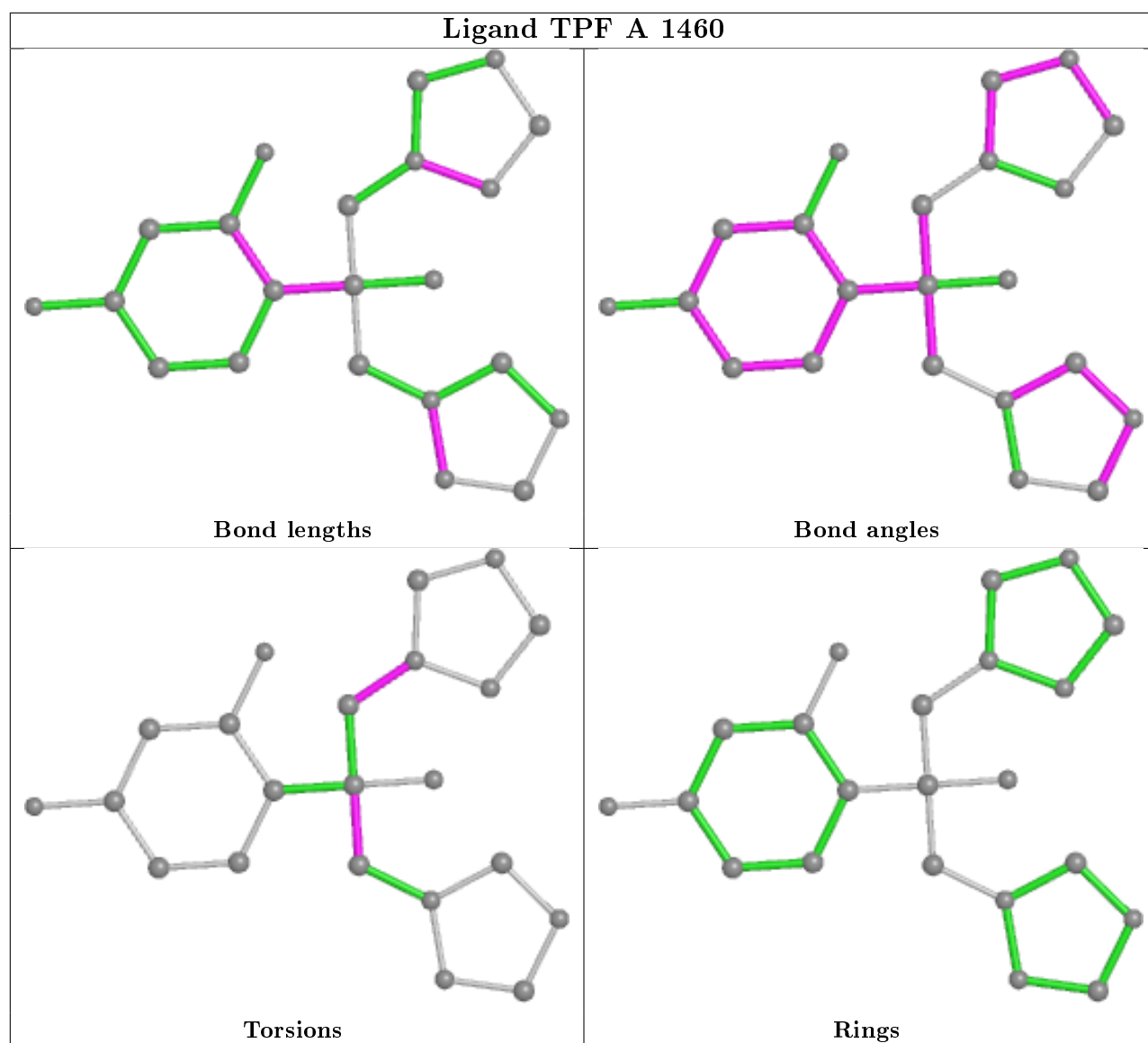
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	1460	TPF	1	0
2	A	1450	HEM	7	0
2	B	1450	HEM	3	0
3	A	1460	TPF	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.









## 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 [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	441/473 (93%)	-0.37	4 (0%) 84 87	8, 22, 46, 56	0
1	B	444/473 (93%)	-0.31	6 (1%) 75 79	9, 22, 43, 51	0
All	All	885/946 (93%)	-0.34	10 (1%) 80 84	8, 22, 45, 56	0

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	245	VAL	3.0
1	A	477	LYS	2.8
1	A	245	VAL	2.7
1	B	196	HIS	2.6
1	A	248	GLU	2.5
1	B	484	HIS	2.2
1	B	230	ARG	2.2
1	B	120	TYR	2.2
1	A	202	SER	2.2
1	B	74	GLY	2.1

### 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](#)

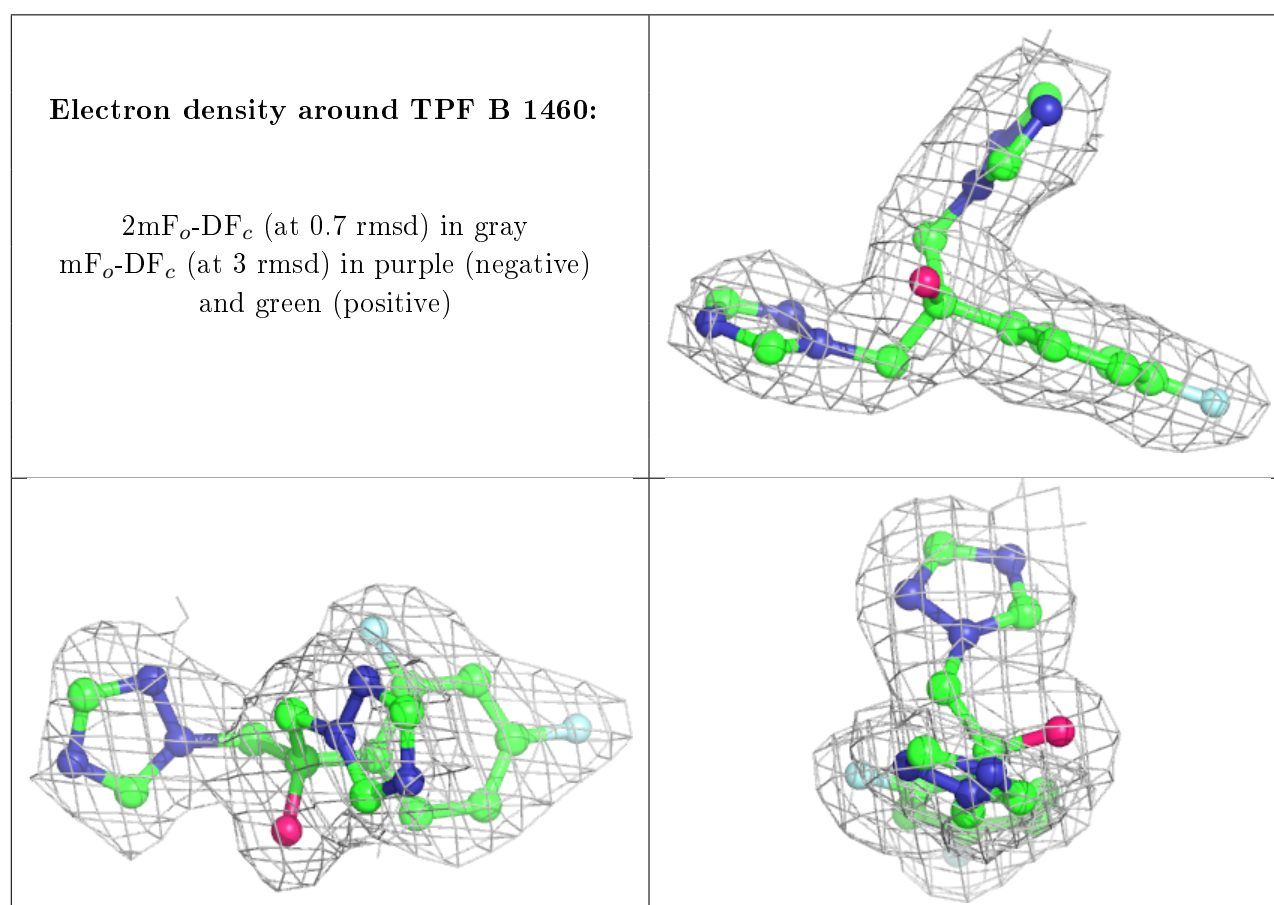
There are no carbohydrates in this entry.

## 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, 95<sup>th</sup> 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	TPF	B	1460	22/22	0.95	0.12	21,32,35,36	0
3	TPF	A	1460	22/22	0.95	0.12	27,32,40,41	0
2	HEM	B	1450	43/43	0.98	0.08	5,12,23,24	0
2	HEM	A	1450	43/43	0.98	0.09	11,16,22,23	0

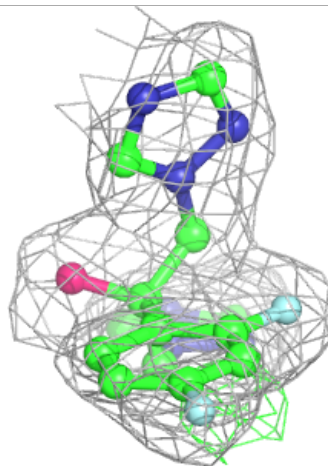
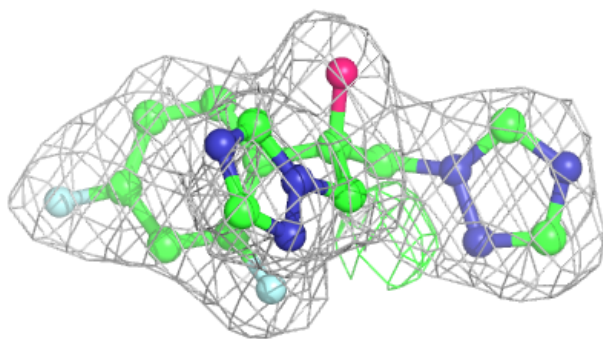
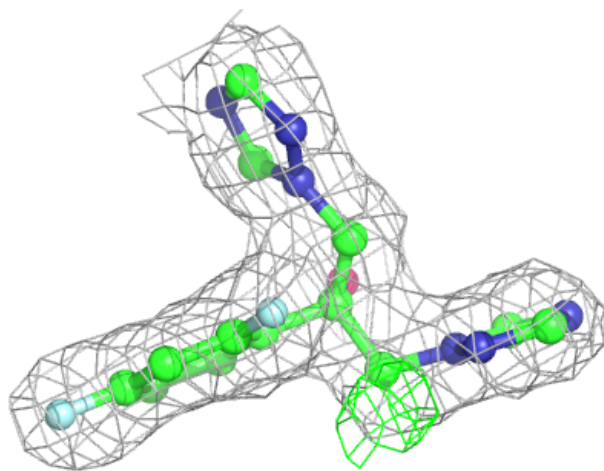
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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





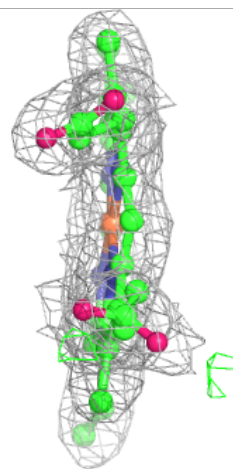
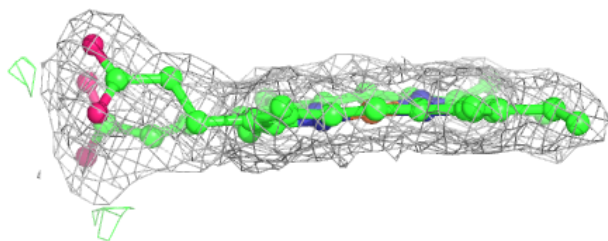
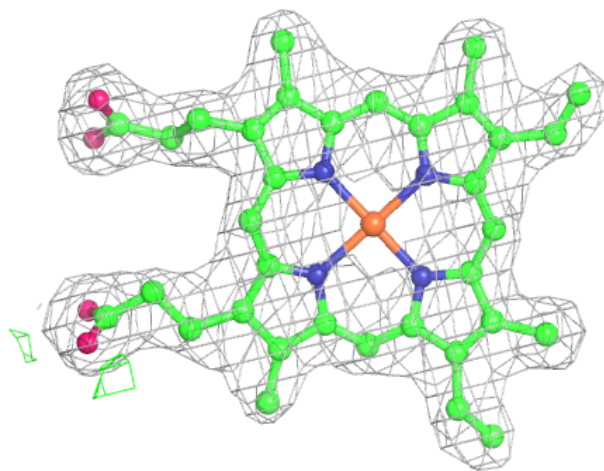
**Electron density around TPF A 1460:**

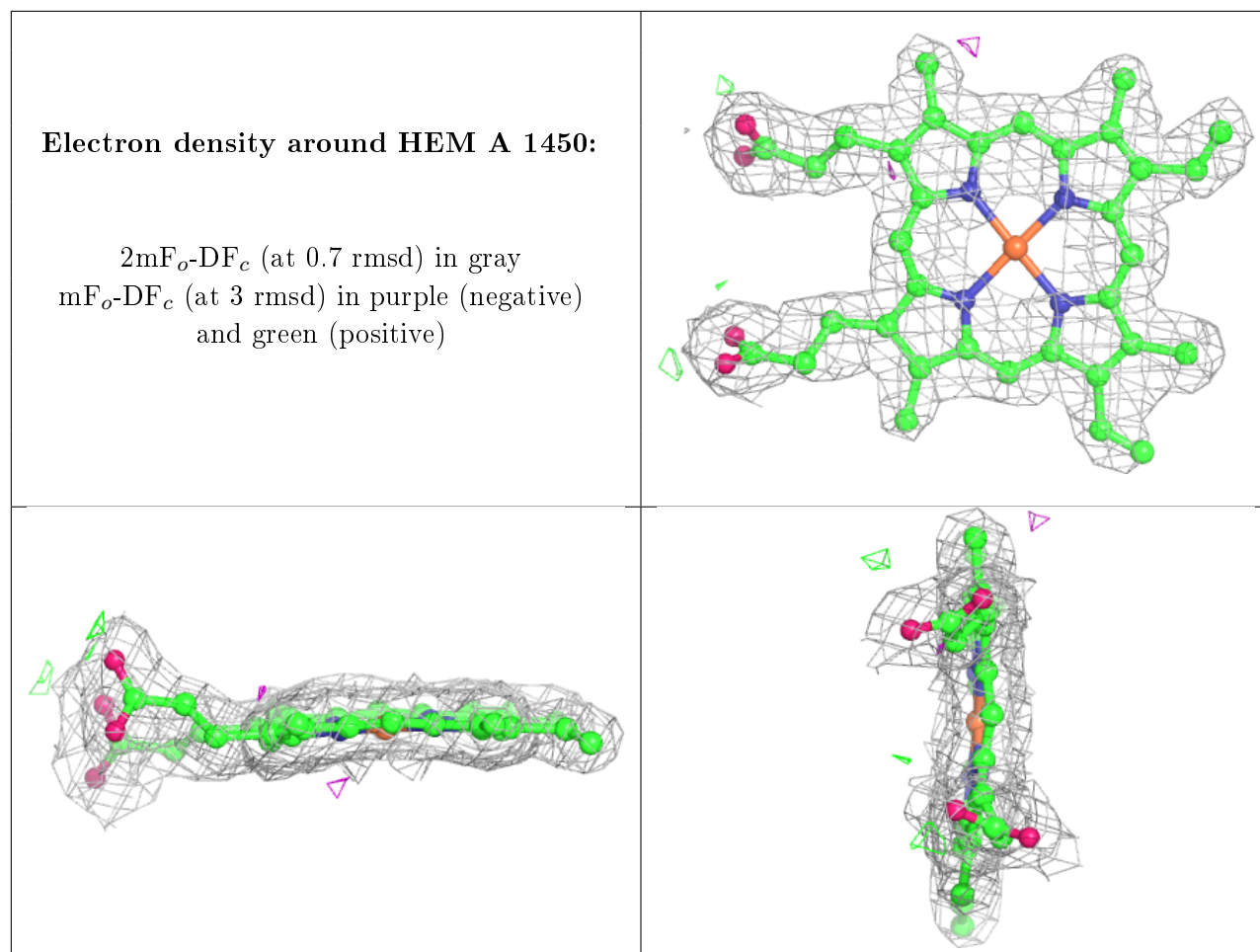
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



**Electron density around HEM B 1450:**

$2mF_o-DF_c$  (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.