



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 06:18 PM UTC

PDB ID : 4KTF / pdb_00004ktf
Title : Crystal structure of Mycobacterium tuberculosis CYP121 in complex with 4, 4'-(3-amino-1H-pyrazole-4,5-diyl)diphenol
Authors : Hudson, S.A.
Deposited on : 2013-05-20
Resolution : 1.35 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

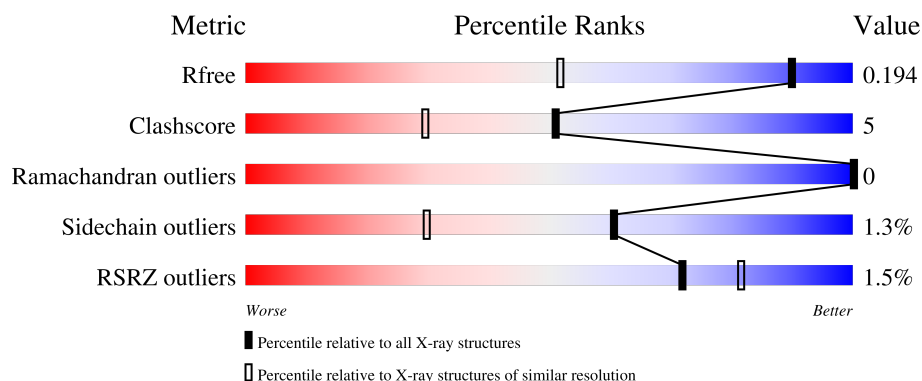
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

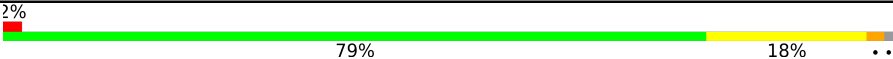
The reported resolution of this entry is 1.35 Å.

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	1216 (1.36-1.36)
Clashscore	190562	1232 (1.36-1.36)
Ramachandran outliers	187476	1220 (1.36-1.36)
Sidechain outliers	187428	1220 (1.36-1.36)
RSRZ outliers	180081	1214 (1.36-1.36)

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	395	

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 3739 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 P450 121.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	393	Total	C	N	O	S	0	25	1
			3068	1963	533	561	11			

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



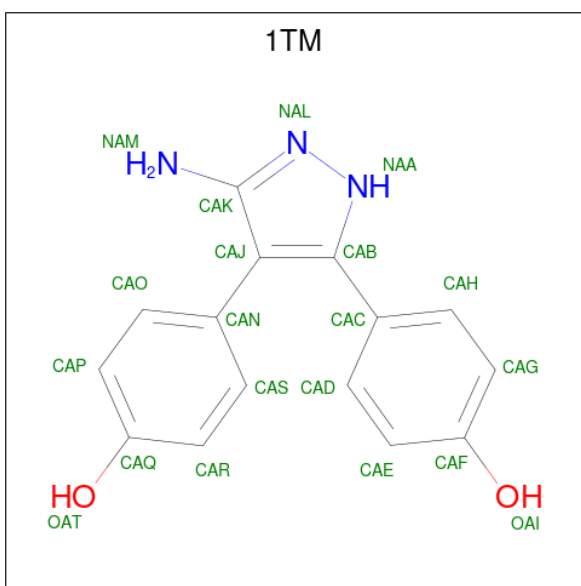
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O_4S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		

- Molecule 4 is 4,4'-(3-amino-1H-pyrazole-4,5-diyl)diphenol (CCD ID: 1TM) (formula: C₁₅H₁₃N₃O₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			20	15	3	2		

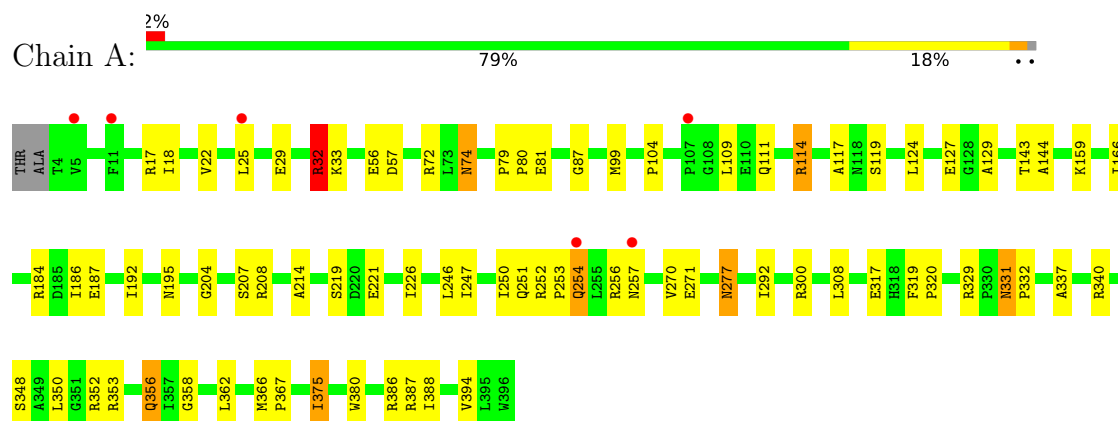
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	588	Total	O	0	0
			588	588		

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: Cytochrome P450 121



4 Data and refinement statistics

Property	Value	Source
Space group	P 65 2 2	Depositor
Cell constants a, b, c, α , β , γ	78.20Å 78.20Å 263.19Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	35.71 – 1.35 35.71 – 1.35	Depositor EDS
% Data completeness (in resolution range)	91.3 (35.71-1.35) 91.3 (35.71-1.35)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.07 (at 1.35Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.166 , 0.195 0.165 , 0.194	Depositor DCC
R_{free} test set	4818 reflections (4.58%)	wwPDB-VP
Wilson B-factor (Å ²)	11.3	Xtriage
Anisotropy	0.039	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 36.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	3739	wwPDB-VP
Average B, all atoms (Å ²)	13.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.77% 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: SO4, HEM, 1TM

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	1.88	48/3244 (1.5%)	1.51	25/4412 (0.6%)

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	0	1

All (48) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	32	ARG	CZ-NH1	8.38	1.44	1.32
1	A	87	GLY	N-CA	-7.53	1.36	1.45
1	A	271[A]	GLU	N-CA	-7.51	1.37	1.46
1	A	271[B]	GLU	N-CA	-7.51	1.37	1.46
1	A	204	GLY	N-CA	7.35	1.54	1.45
1	A	253	PRO	N-CA	-6.81	1.38	1.47
1	A	254	GLN	CA-C	6.64	1.61	1.52
1	A	253	PRO	CA-C	6.42	1.61	1.52
1	A	319	PHE	CA-C	6.38	1.59	1.52
1	A	117	ALA	CA-CB	-6.29	1.43	1.53
1	A	388	ILE	CA-CB	-6.23	1.46	1.54
1	A	219	SER	N-CA	-6.22	1.38	1.45
1	A	308	LEU	C-O	6.12	1.31	1.23
1	A	32	ARG	CA-CB	-6.09	1.43	1.53
1	A	18	ILE	CA-C	6.06	1.59	1.52
1	A	319	PHE	N-CA	-5.97	1.39	1.46
1	A	367	PRO	N-CA	-5.95	1.39	1.47
1	A	144	ALA	C-O	5.85	1.30	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	337	ALA	N-CA	5.85	1.53	1.46
1	A	79	PRO	CA-C	-5.83	1.48	1.51
1	A	124	LEU	N-CA	5.79	1.53	1.46
1	A	207[A]	SER	N-CA	-5.77	1.39	1.46
1	A	207[B]	SER	N-CA	-5.77	1.39	1.46
1	A	57	ASP	C-O	5.76	1.30	1.24
1	A	356	GLN	N-CA	5.69	1.53	1.46
1	A	256	ARG	CZ-NH2	5.61	1.40	1.33
1	A	292	ILE	C-O	5.54	1.29	1.24
1	A	81	GLU	CD-OE2	5.50	1.35	1.25
1	A	332	PRO	C-O	5.46	1.30	1.24
1	A	394	VAL	N-CA	5.43	1.52	1.46
1	A	356	GLN	CA-C	-5.30	1.46	1.52
1	A	25	LEU	CA-CB	5.26	1.61	1.53
1	A	114[A]	ARG	N-CA	-5.25	1.40	1.46
1	A	114[B]	ARG	N-CA	-5.25	1.40	1.46
1	A	350	LEU	N-CA	-5.21	1.40	1.46
1	A	214	ALA	C-O	5.19	1.30	1.24
1	A	320	PRO	N-CA	-5.17	1.40	1.47
1	A	387	ARG	CD-NE	-5.15	1.39	1.46
1	A	18	ILE	C-O	-5.15	1.18	1.24
1	A	277	ASN	N-CA	5.12	1.52	1.46
1	A	195	ASN	C-N	5.11	1.40	1.33
1	A	184	ARG	C-O	5.11	1.30	1.24
1	A	375	ILE	CB-CG1	5.10	1.63	1.53
1	A	186	ILE	C-N	-5.09	1.27	1.33
1	A	192	ILE	C-O	5.05	1.30	1.24
1	A	22	VAL	CA-C	5.03	1.58	1.52
1	A	380	TRP	N-CA	5.01	1.52	1.46
1	A	129	ALA	N-CA	5.00	1.52	1.45

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	32	ARG	NE-CZ-NH1	8.63	130.13	121.50
1	A	32	ARG	NE-CZ-NH2	-8.33	111.71	119.20
1	A	250	ILE	CA-C-O	-7.15	112.92	120.57
1	A	252	ARG	O-C-N	-6.79	115.29	121.34
1	A	352	ARG	NE-CZ-NH1	-6.46	115.04	121.50
1	A	366	MET	O-C-N	-6.34	115.88	121.20
1	A	252	ARG	CA-C-O	6.32	129.11	121.02
1	A	308	LEU	CA-C-O	-6.30	114.17	120.98

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	348	SER	N-CA-CB	5.96	118.66	110.01
1	A	221	GLU	CB-CG-CD	5.94	122.69	112.60
1	A	208	ARG	NE-CZ-NH1	-5.85	115.65	121.50
1	A	329	ARG	NE-CZ-NH2	5.79	124.41	119.20
1	A	99[A]	MET	CB-CA-C	-5.59	101.51	110.79
1	A	99[B]	MET	CB-CA-C	-5.59	101.51	110.79
1	A	254	GLN	CA-C-O	5.58	126.47	120.55
1	A	358	GLY	O-C-N	5.50	127.47	122.19
1	A	253	PRO	N-CA-C	5.49	121.92	113.75
1	A	143	THR	N-CA-C	-5.41	105.30	111.14
1	A	219	SER	N-CA-C	5.41	118.12	110.50
1	A	226	ILE	CA-C-N	5.30	125.98	119.99
1	A	226	ILE	C-N-CA	5.30	125.98	119.99
1	A	270	VAL	O-C-N	-5.26	116.56	121.87
1	A	256	ARG	NE-CZ-NH1	-5.24	116.26	121.50
1	A	329	ARG	NH1-CZ-NH2	-5.11	112.66	119.30
1	A	166	ILE	N-CA-C	-5.06	107.80	111.90

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	72	ARG	Sidechain

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	3068	0	3083	32	0
2	A	43	0	30	1	0
3	A	20	0	0	0	0
4	A	20	0	12	0	0
5	A	588	0	0	13	2
All	All	3739	0	3125	33	2

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 5.

All (33) 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:17:ARG:NE	5:A:1042:HOH:O	2.12	0.82
1:A:251[B]:GLN:NE2	1:A:375:ILE:HD13	1.95	0.81
1:A:353:ARG:HH11	1:A:356:GLN:HE21	1.26	0.80
1:A:251[B]:GLN:NE2	1:A:375:ILE:CD1	2.45	0.79
1:A:127[A]:GLU:CG	5:A:1043:HOH:O	2.34	0.75
1:A:127[B]:GLU:CG	5:A:1043:HOH:O	2.38	0.71
1:A:29:GLU:HG2	1:A:32:ARG:HB3	1.73	0.70
1:A:187[A]:GLU:OE2	5:A:1027:HOH:O	2.11	0.69
1:A:254:GLN:CG	5:A:981:HOH:O	2.41	0.68
1:A:74:ASN:C	1:A:74:ASN:HD22	2.02	0.67
1:A:254:GLN:HG3	5:A:981:HOH:O	1.99	0.62
1:A:353:ARG:HH11	1:A:356:GLN:NE2	1.96	0.61
1:A:300:ARG:CG	5:A:1079:HOH:O	2.48	0.60
1:A:111:GLN:NE2	1:A:114[B]:ARG:HH21	1.98	0.60
1:A:254:GLN:HG2	5:A:981:HOH:O	2.03	0.57
1:A:247:ILE:O	1:A:251[B]:GLN:HG3	2.09	0.53
1:A:331:ASN:H	1:A:331:ASN:HD22	1.56	0.52
2:A:401:HEM:HMC1	2:A:401:HEM:HBC2	1.93	0.51
1:A:257:ASN:ND2	5:A:1053:HOH:O	2.22	0.49
1:A:29:GLU:CG	1:A:32:ARG:HB3	2.43	0.48
1:A:317:GLU:HG3	5:A:1066:HOH:O	2.12	0.48
1:A:159[B]:LYS:CG	5:A:1026:HOH:O	2.63	0.46
1:A:56:GLU:OE2	1:A:340[B]:ARG:HD2	2.16	0.46
1:A:246:LEU:HD22	1:A:362[B]:LEU:HD22	1.99	0.45
1:A:104:PRO:HA	1:A:109:LEU:HD23	1.98	0.44
1:A:251[B]:GLN:HE22	1:A:375:ILE:HD13	1.76	0.44
1:A:331:ASN:HD22	1:A:331:ASN:N	2.16	0.41
1:A:33[B]:LYS:CD	5:A:877:HOH:O	2.68	0.41
1:A:80[A]:PRO:HG3	5:A:947:HOH:O	2.21	0.41
1:A:246:LEU:HD22	1:A:362[B]:LEU:CD2	2.50	0.41

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:A:849:HOH:O	5:A:849:HOH:O[10_665]	1.48	0.72
5:A:814:HOH:O	5:A:953:HOH:O[10_555]	1.96	0.24

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	417/395 (106%)	409 (98%)	8 (2%)	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	330/325 (102%)	326 (99%)	4 (1%)	63	32

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	32	ARG
1	A	74	ASN
1	A	277	ASN
1	A	331	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	74	ASN
1	A	111	GLN
1	A	154	GLN
1	A	331	ASN

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Mol	Chain	Res	Type
1	A	356	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](#)

6 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	SO4	A	403	-	4,4,4	0.73	0	6,6,6	0.57	0
3	SO4	A	405	-	4,4,4	1.05	0	6,6,6	0.81	0
3	SO4	A	402	-	4,4,4	1.66	2 (50%)	6,6,6	2.00	1 (16%)
2	HEM	A	401	5,1	50,50,50	1.70	12 (24%)	67,82,82	1.86	11 (16%)
4	1TM	A	406	-	22,22,22	2.00	8 (36%)	26,31,31	2.14	12 (46%)
3	SO4	A	404	-	4,4,4	1.03	0	6,6,6	2.28	2 (33%)

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
4	1TM	A	406	-	-	0/8/8/8	0/3/3/3
2	HEM	A	401	5,1	-	0/14/54/54	-

All (22) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	401	HEM	FE-NB	4.69	2.09	1.94
2	A	401	HEM	C4B-NB	-3.98	1.31	1.38
4	A	406	1TM	CAK-NAM	3.93	1.44	1.34
2	A	401	HEM	C1B-NB	-3.88	1.33	1.40
4	A	406	1TM	CAS-CAR	3.60	1.44	1.38
4	A	406	1TM	CAP-CAO	3.22	1.44	1.38
2	A	401	HEM	CHB-C4A	-3.01	1.32	1.39
4	A	406	1TM	CAC-CAB	-2.71	1.43	1.47
2	A	401	HEM	C4D-ND	-2.55	1.35	1.40
3	A	402	SO4	O2-S	2.45	1.59	1.44
2	A	401	HEM	FE-NA	2.42	2.03	1.95
4	A	406	1TM	CAK-NAL	2.42	1.35	1.33
4	A	406	1TM	CAN-CAJ	-2.41	1.45	1.49
4	A	406	1TM	CAG-CAF	2.36	1.43	1.39
2	A	401	HEM	C1D-C2D	2.32	1.49	1.44
4	A	406	1TM	OAT-CAQ	2.31	1.42	1.37
2	A	401	HEM	O1D-CGD	2.28	1.29	1.22
2	A	401	HEM	C4D-C3D	2.24	1.48	1.45
2	A	401	HEM	FE-ND	2.23	2.01	1.94
3	A	402	SO4	O1-S	2.21	1.58	1.44
2	A	401	HEM	CHB-C1B	2.01	1.42	1.38
2	A	401	HEM	O1A-CGA	2.00	1.28	1.22

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	401	HEM	C1B-NB-C4B	7.88	114.54	105.21
2	A	401	HEM	C3B-C4B-NB	-6.35	104.91	109.47
2	A	401	HEM	CHC-C4B-NB	5.44	130.27	124.42
4	A	406	1TM	CAC-CAB-CAJ	4.78	138.78	131.46
3	A	404	SO4	O4-S-O3	4.18	131.59	108.54
4	A	406	1TM	CAC-CAB-NAA	-4.07	116.08	121.23
2	A	401	HEM	C4C-C3C-C2C	3.60	109.94	106.81
4	A	406	1TM	CAR-CAS-CAN	3.49	124.52	120.80
3	A	404	SO4	O4-S-O2	-3.45	91.54	109.56
3	A	402	SO4	O4-S-O3	3.40	127.28	108.54
2	A	401	HEM	C2B-C1B-NB	-3.37	105.96	109.84

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	406	1TM	CAS-CAN-CAJ	2.85	125.27	120.79
2	A	401	HEM	CHB-C1B-NB	2.83	127.87	124.37
4	A	406	1TM	CAG-CAH-CAC	2.75	123.73	120.80
4	A	406	1TM	CAN-CAJ-CAB	2.69	132.18	127.63
2	A	401	HEM	CHD-C4C-C3C	2.67	129.71	125.21
4	A	406	1TM	CAH-CAC-CAB	2.64	123.90	120.76
4	A	406	1TM	CAS-CAN-CAO	-2.61	115.25	118.57
4	A	406	1TM	CAH-CAC-CAD	-2.26	115.70	118.57
4	A	406	1TM	CAB-NAA-NAL	2.23	116.66	112.88
4	A	406	1TM	CAE-CAD-CAC	2.23	123.18	120.80
2	A	401	HEM	O2D-CGD-CBD	2.19	120.90	114.00
4	A	406	1TM	CAS-CAR-CAQ	-2.15	117.60	119.88
2	A	401	HEM	O2A-CGA-CBA	2.10	120.63	114.00
2	A	401	HEM	C3C-C2C-C1C	-2.05	105.11	107.05
2	A	401	HEM	O2A-CGA-O1A	-2.02	118.13	123.33

There are no chirality outliers.

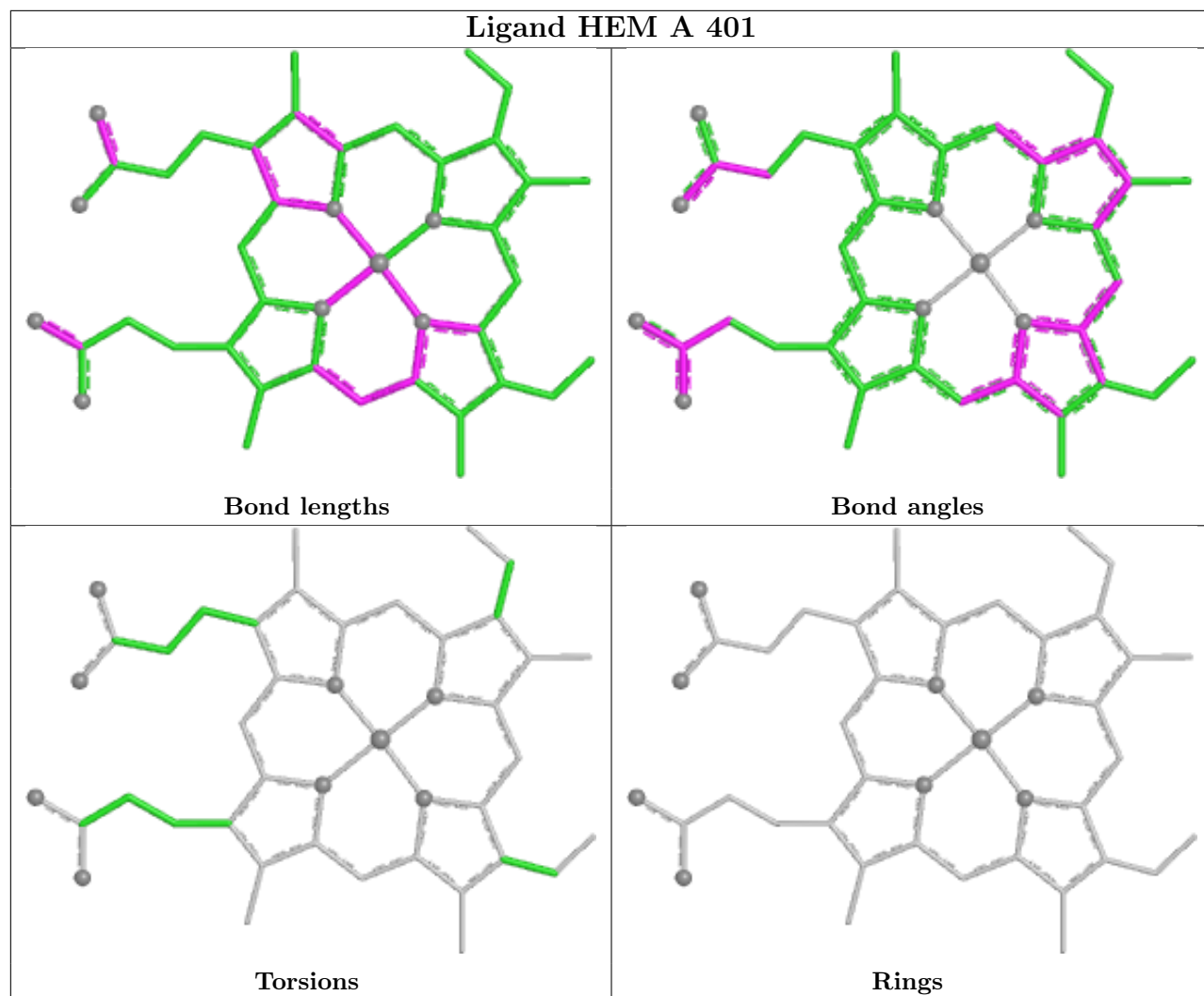
There are no torsion outliers.

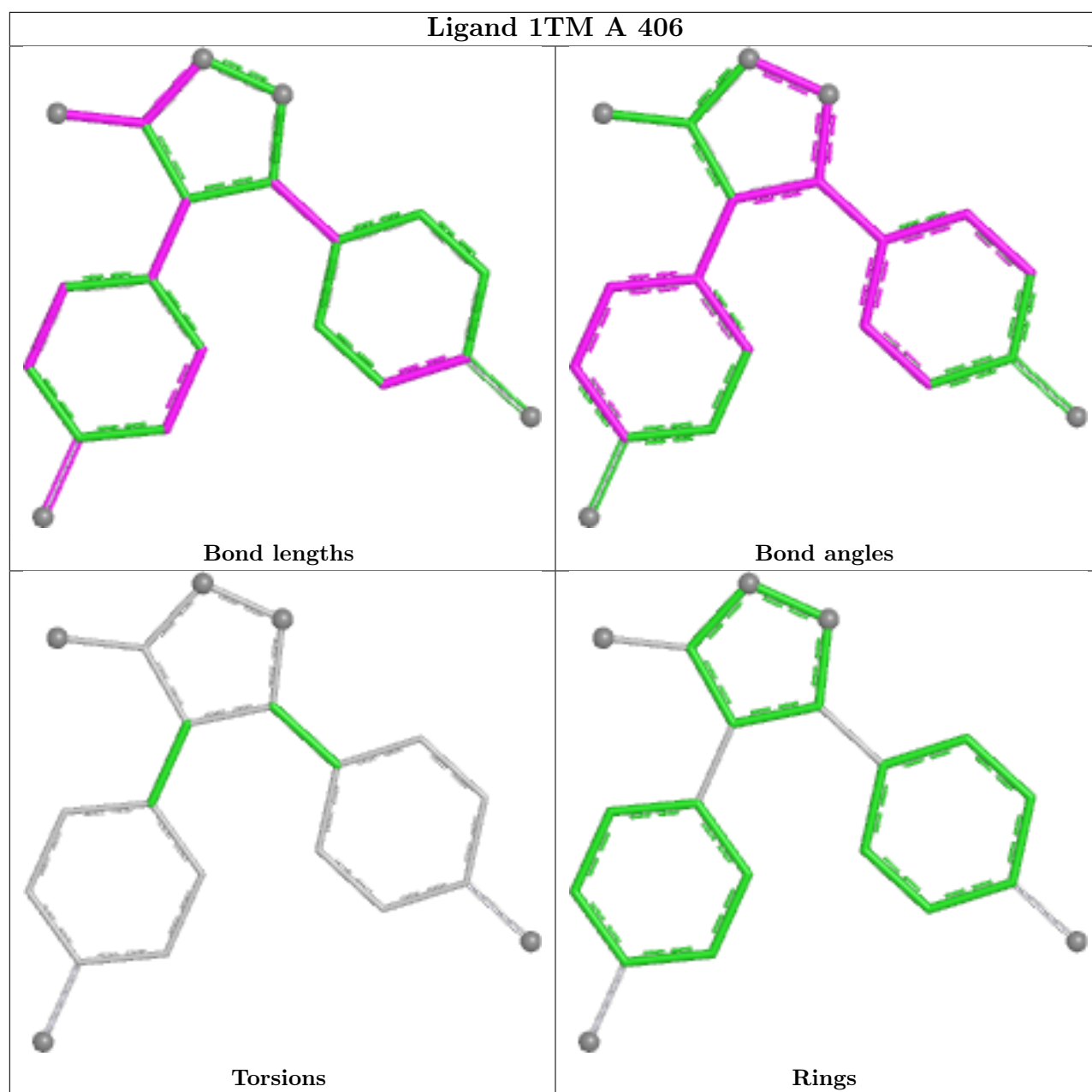
There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	401	HEM	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, 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	393/395 (99%)	-0.06	6 (1%) 72 81	4, 10, 22, 33	26 (6%)

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	5	VAL	3.5
1	A	107	PRO	3.3
1	A	254	GLN	2.5
1	A	257	ASN	2.3
1	A	25	LEU	2.1
1	A	11	PHE	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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

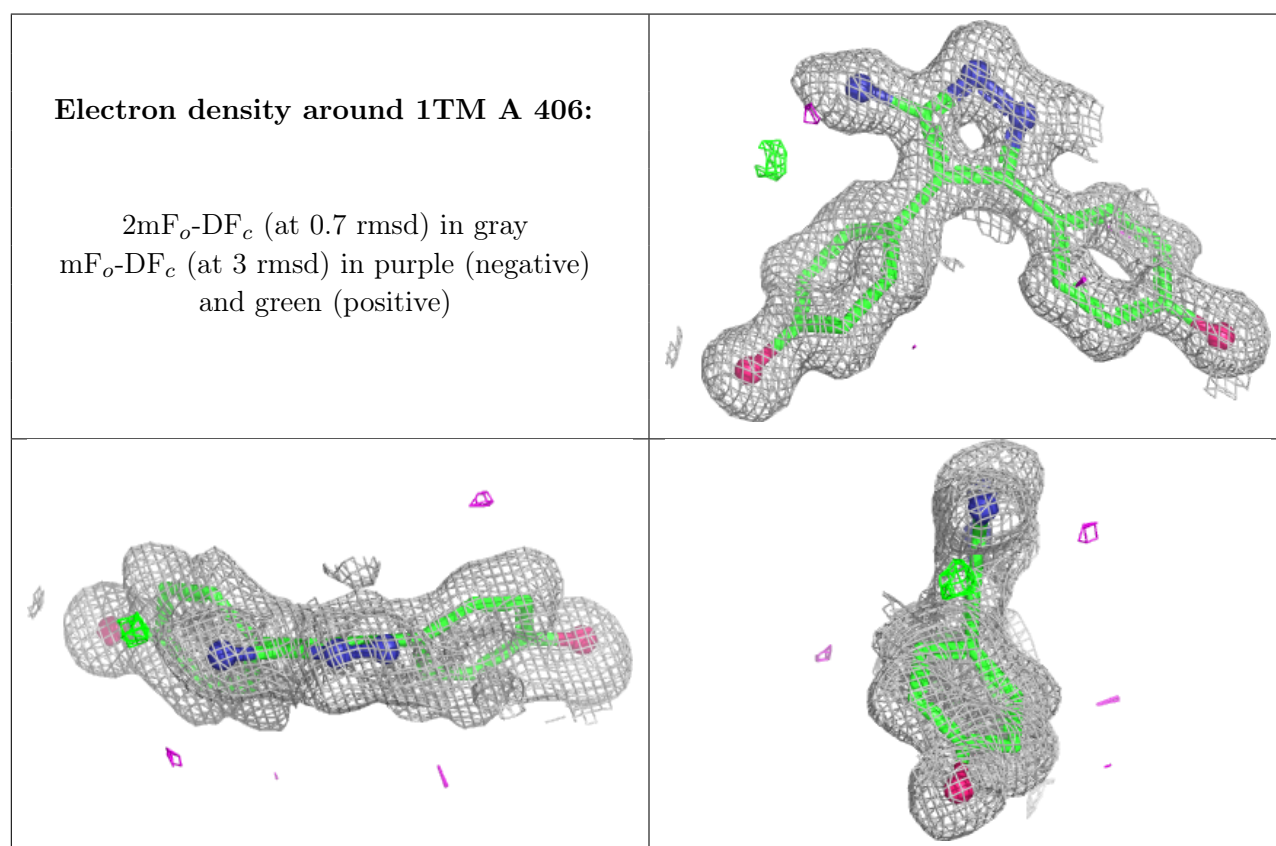
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	SO4	A	403	5/5	0.91	0.10	27,33,38,40	0

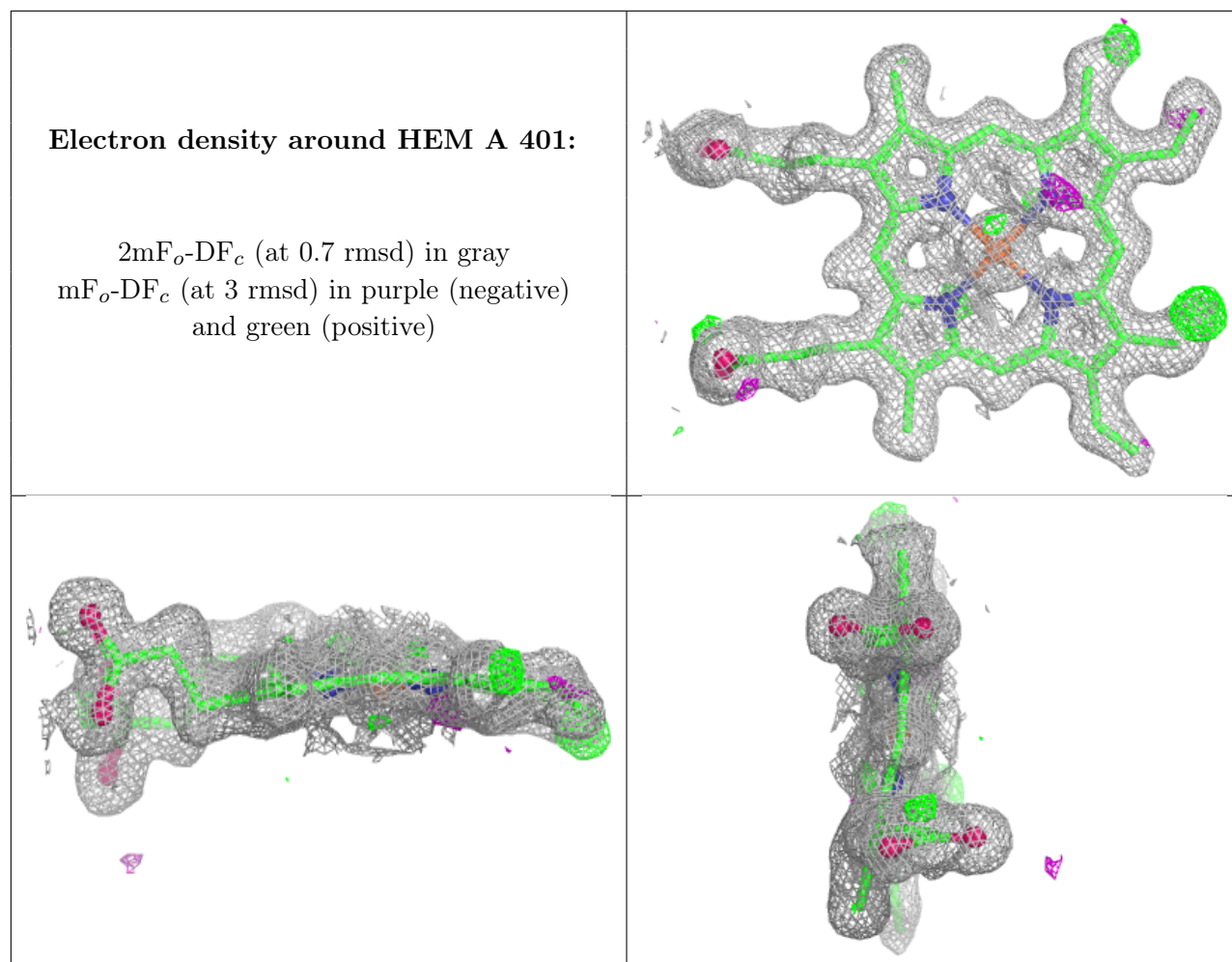
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	SO4	A	405	5/5	0.93	0.11	16,25,32,32	0
3	SO4	A	402	5/5	0.94	0.13	21,23,27,34	0
3	SO4	A	404	5/5	0.97	0.08	15,16,22,25	0
4	1TM	A	406	20/20	0.98	0.04	6,8,10,10	0
2	HEM	A	401	43/43	0.99	0.05	5,8,11,13	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.





6.5 Other polymers ⓘ

There are no such residues in this entry.