



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 18, 2026 – 01:58 PM UTC

PDB ID : 3C8Z / pdb_00003c8z
Title : The 1.6 Å Crystal Structure of MshC: The Rate Limiting Enzyme in the
Mycothioli Biosynthetic Pathway
Authors : Tremblay, L.W.; Fan, F.; Vetting, M.W.; Blanchard, J.S.
Deposited on : 2008-02-14
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)	:	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

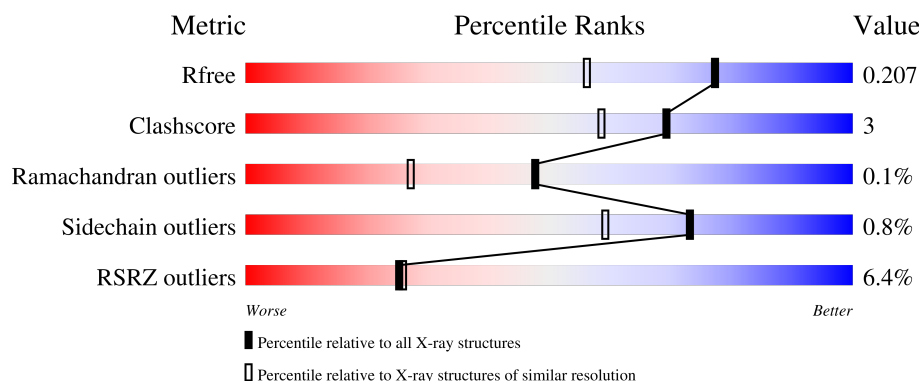
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.60 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))
R_{free}	180053	4673 (1.60-1.60)
Clashscore	190562	4931 (1.60-1.60)
Ramachandran outliers	187476	4831 (1.60-1.60)
Sidechain outliers	187428	4830 (1.60-1.60)
RSRZ outliers	180081	4672 (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 $\leq 5\%$. 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	414	7% 90% 7% .
1	B	414	6% 90% 7% .

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
4	5CA	A	418	X	-	-	-
4	5CA	B	416	X	-	-	-

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 7207 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 CysteinyI-tRNA synthetase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	399	Total	C	N	O	S	0	0	0
			3100	1948	550	593	9			
1	B	401	Total	C	N	O	S	0	0	0
			3118	1959	554	595	10			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	HIS	-	expression tag	UNP A0QZY0
A	0	MET	-	expression tag	UNP A0QZY0
B	-1	HIS	-	expression tag	UNP A0QZY0
B	0	MET	-	expression tag	UNP A0QZY0

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

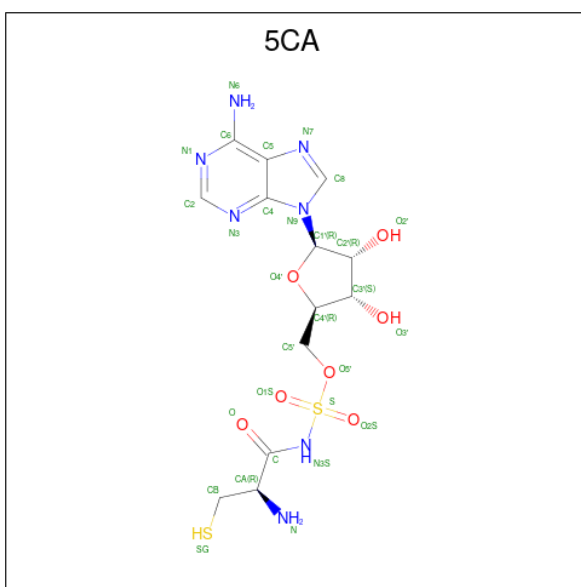
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Zn	0	0
			1	1		
2	B	1	Total	Zn	0	0
			1	1		

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



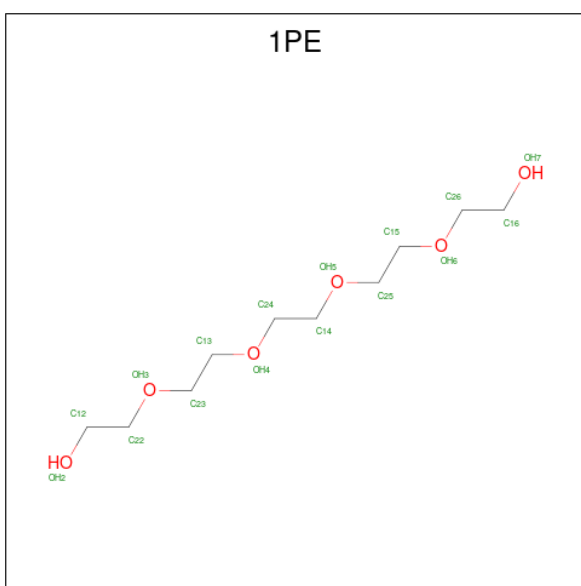
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 4 is 5'-O-(N-(L-CYSTEINYL)-SULFAMOYL)ADENOSINE (CCD ID: 5CA) (formula: C₁₃H₁₉N₇O₇S₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total 29	C 13	N 7	O 7	S 2	0	0
4	B	1	Total 29	C 13	N 7	O 7	S 2	0	0

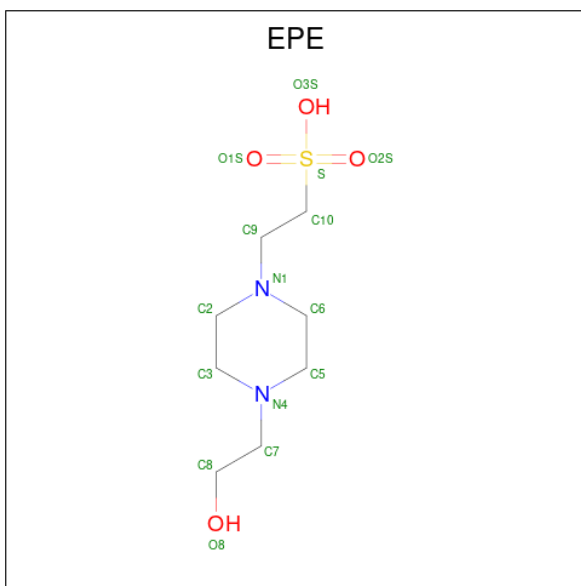
- Molecule 5 is PENTAETHYLENE GLYCOL (CCD ID: 1PE) (formula: $C_{10}H_{22}O_6$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			16	10	6		

- Molecule 6 is 4-(2-HYDROXYETHYL)-1-PIPERAZINE ETHANESULFONIC ACID (CCD

ID: EPE) (formula: C₈H₁₈N₂O₄S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	B	1	Total	C	N	O	S	0	0
			15	8	2	4	1		

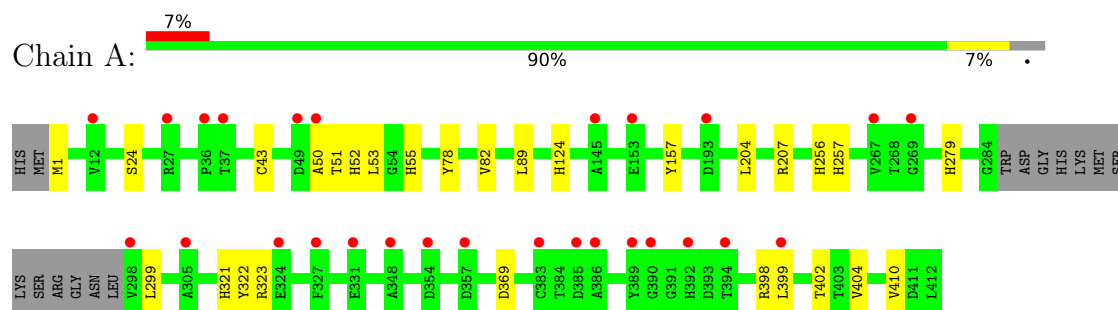
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	396	Total	O	0	0
			396	396		
7	B	472	Total	O	0	0
			472	472		

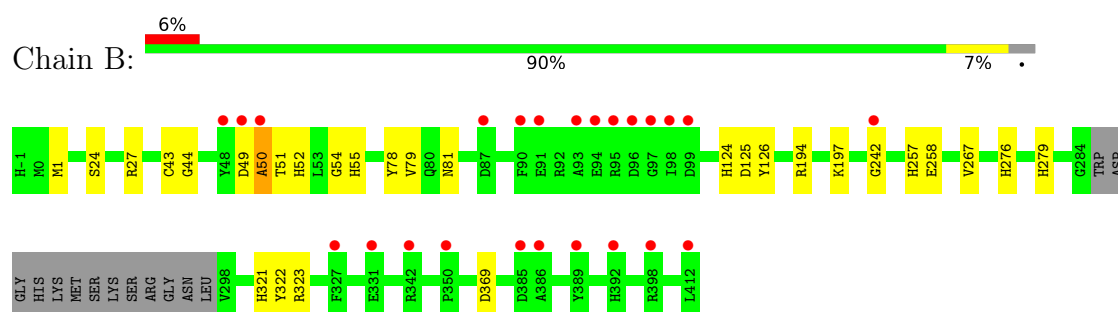
3 Residue-property plots [i](#)

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 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: Cysteinyl-tRNA synthetase



• Molecule 1: Cysteinyl-tRNA synthetase



4 Data and refinement statistics

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, α , β , γ	123.52Å 123.52Å 186.03Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	37.09 – 1.60 37.09 – 1.60	Depositor EDS
% Data completeness (in resolution range)	99.7 (37.09-1.60) 99.7 (37.09-1.60)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.85 (at 1.60Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.185 , 0.207 0.184 , 0.207	Depositor DCC
R_{free} test set	6973 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	20.7	Xtriage
Anisotropy	0.016	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 40.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.017 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	7207	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.11% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, EPE, 1PE, SO4, 5CA

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	0.64	2/3181 (0.1%)	0.78	0/4344
1	B	0.64	0/3200	0.80	2/4369 (0.0%)
All	All	0.64	2/6381 (0.0%)	0.79	2/8713 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	51	THR	C-O	-5.67	1.16	1.24
1	A	256	HIS	C-O	-5.04	1.18	1.24

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	258	GLU	N-CA-C	7.40	119.99	111.11
1	B	242	GLY	N-CA-C	6.72	121.86	114.40

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	3100	0	2957	19	0
1	B	3118	0	2973	24	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	20	0	0	0	0
3	B	10	0	0	0	0
4	A	29	0	14	1	0
4	B	29	0	16	1	0
5	A	16	0	22	1	0
6	B	15	0	18	0	0
7	A	396	0	0	0	0
7	B	472	0	0	1	0
All	All	7207	0	6000	43	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 3.

All (43) 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:50:ALA:HB1	1:B:51:THR:HA	1.26	1.10
1:B:24:SER:HB3	7:B:840:HOH:O	1.77	0.84
1:B:50:ALA:HB1	1:B:51:THR:CA	2.07	0.84
1:A:78:TYR:H	1:A:124:HIS:HD2	1.24	0.83
1:B:78:TYR:H	1:B:124:HIS:HD2	1.26	0.83
1:B:50:ALA:CB	1:B:51:THR:HA	2.11	0.78
1:A:257:HIS:ND1	1:A:279:HIS:HE1	1.81	0.77
1:B:78:TYR:H	1:B:124:HIS:CD2	2.05	0.74
1:B:257:HIS:ND1	1:B:279:HIS:HE1	1.88	0.71
1:A:24:SER:OG	1:A:279:HIS:HD2	1.74	0.69
1:B:24:SER:OG	1:B:279:HIS:HD2	1.76	0.69
1:A:78:TYR:H	1:A:124:HIS:CD2	2.08	0.67
1:B:50:ALA:HB1	1:B:55:HIS:HD2	1.59	0.67
1:A:52:HIS:H	1:A:55:HIS:HD2	1.44	0.66
1:B:50:ALA:CB	1:B:55:HIS:CD2	2.82	0.62
1:B:50:ALA:HB1	1:B:55:HIS:CD2	2.37	0.60
1:B:49:ASP:O	1:B:50:ALA:HB3	2.04	0.58
1:A:321:HIS:HE1	1:A:369:ASP:O	1.91	0.53
1:B:321:HIS:HE1	1:B:369:ASP:O	1.91	0.53
1:A:43:CYS:HB3	4:A:418:5CA:HB2	1.91	0.53
1:B:1:MET:HE3	1:B:126:TYR:CD2	2.44	0.53
1:A:157:TYR:CE1	5:A:419:1PE:H162	2.44	0.52
1:A:52:HIS:H	1:A:55:HIS:CD2	2.25	0.52

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:257:HIS:ND1	1:A:279:HIS:CE1	2.71	0.52
1:B:50:ALA:HB2	1:B:55:HIS:CD2	2.47	0.50
1:B:43:CYS:HB3	4:B:416:5CA:HB2	1.94	0.50
1:B:52:HIS:H	1:B:55:HIS:HD2	1.62	0.48
1:B:194:ARG:HB3	1:B:197:LYS:HD3	1.95	0.48
1:A:404:VAL:HG12	1:A:410:VAL:HB	1.97	0.47
1:A:398:ARG:O	1:A:402:THR:HG23	2.14	0.47
1:A:50:ALA:HB3	1:A:299:LEU:HD22	1.97	0.47
1:A:404:VAL:CG1	1:A:410:VAL:HB	2.45	0.46
1:A:322:TYR:CE2	1:A:323:ARG:HG3	2.52	0.45
1:B:44:GLY:HA3	1:B:81:ASN:O	2.18	0.44
1:B:54:GLY:O	1:B:55:HIS:C	2.61	0.44
1:A:50:ALA:HB3	1:A:299:LEU:HD13	2.00	0.43
1:B:52:HIS:H	1:B:55:HIS:CD2	2.37	0.43
1:B:257:HIS:ND1	1:B:279:HIS:CE1	2.78	0.43
1:A:1:MET:HE1	1:A:82:VAL:HG21	1.99	0.43
1:A:204:LEU:HD21	1:A:207:ARG:HB2	2.01	0.42
1:B:79:VAL:HG22	1:B:125:ASP:HB2	2.01	0.42
1:A:53:LEU:HD23	1:A:53:LEU:HA	1.84	0.41
1:B:322:TYR:CE2	1:B:323:ARG:HG3	2.56	0.41

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	395/414 (95%)	392 (99%)	3 (1%)	0	100	100
1	B	397/414 (96%)	390 (98%)	6 (2%)	1 (0%)	36	20
All	All	792/828 (96%)	782 (99%)	9 (1%)	1 (0%)	48	27

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	50	ALA

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	317/330 (96%)	315 (99%)	2 (1%)	78	66
1	B	319/330 (97%)	316 (99%)	3 (1%)	70	55
All	All	636/660 (96%)	631 (99%)	5 (1%)	73	59

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	89	LEU
1	A	399	LEU
1	B	27	ARG
1	B	267	VAL
1	B	276	HIS

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	55	HIS
1	A	80	GLN
1	A	124	HIS
1	A	263	HIS
1	A	279	HIS
1	A	306	GLN
1	A	321	HIS
1	B	55	HIS
1	B	80	GLN
1	B	124	HIS
1	B	246	GLN
1	B	263	HIS
1	B	279	HIS

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	321	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 12 ligands modelled in this entry, 2 are monoatomic - leaving 10 for Mogul analysis.

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	B	415	-	4,4,4	0.25	0	6,6,6	0.09	0
3	SO4	B	414	-	4,4,4	0.33	0	6,6,6	0.36	0
3	SO4	A	414	-	4,4,4	0.23	0	6,6,6	0.09	0
6	EPE	B	417	-	15,15,15	6.14	3 (20%)	19,20,20	1.13	2 (10%)
3	SO4	A	416	-	4,4,4	0.27	0	6,6,6	0.16	0
5	1PE	A	419	-	15,15,15	0.49	0	14,14,14	0.65	0
4	5CA	B	416	2	31,31,31	2.88	14 (45%)	42,46,46	2.62	19 (45%)
3	SO4	A	415	-	4,4,4	0.26	0	6,6,6	0.22	0
4	5CA	A	418	2	31,31,31	3.27	15 (48%)	42,46,46	3.05	23 (54%)
3	SO4	A	417	-	4,4,4	0.26	0	6,6,6	0.10	0

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
5	1PE	A	419	-	-	5/13/13/13	-
4	5CA	A	418	2	2/2/7/8	6/20/37/37	0/3/3/3
6	EPE	B	417	-	-	2/9/19/19	0/1/1/1
4	5CA	B	416	2	1/1/7/8	3/20/37/37	0/3/3/3

All (32) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	B	417	EPE	C10-S	-22.84	1.45	1.77
4	A	418	5CA	O2S-S	8.25	1.49	1.42
4	B	416	5CA	C4-N3	7.20	1.47	1.34
4	B	416	5CA	O1S-S	6.72	1.48	1.42
4	A	418	5CA	O1S-S	6.11	1.47	1.42
4	B	416	5CA	O2S-S	5.96	1.47	1.42
4	A	418	5CA	C4-N3	5.96	1.45	1.34
4	B	416	5CA	O5'-S	-5.52	1.48	1.60
4	A	418	5CA	O5'-S	-5.17	1.49	1.60
4	A	418	5CA	C2-N1	4.96	1.42	1.33
6	B	417	EPE	O3S-S	4.41	1.63	1.47
4	B	416	5CA	C2-N1	4.05	1.41	1.33
4	A	418	5CA	O3'-C3'	-4.04	1.33	1.43
4	A	418	5CA	C2-N3	3.97	1.41	1.33
6	B	417	EPE	O1S-S	3.80	1.55	1.45
4	A	418	5CA	O4'-C1'	3.65	1.50	1.42
4	A	418	5CA	CB-CA	-3.63	1.49	1.53
4	A	418	5CA	C-N3S	-3.61	1.31	1.37
4	A	418	5CA	C5-C6	3.59	1.51	1.41
4	A	418	5CA	C2'-C3'	-3.43	1.44	1.53
4	B	416	5CA	C5-C4	3.40	1.45	1.39
4	B	416	5CA	C2-N3	3.24	1.39	1.33
4	B	416	5CA	C-N3S	-2.90	1.32	1.37
4	B	416	5CA	C5-C6	2.87	1.49	1.41
4	A	418	5CA	O4'-C4'	-2.80	1.38	1.45
4	B	416	5CA	C1'-N9	-2.62	1.39	1.46
4	B	416	5CA	C3'-C4'	-2.49	1.46	1.53
4	A	418	5CA	C8-N7	2.44	1.36	1.31
4	A	418	5CA	C5-C4	2.36	1.43	1.39
4	B	416	5CA	C2'-C3'	-2.27	1.47	1.53
4	B	416	5CA	C5-N7	2.06	1.42	1.39

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	416	5CA	C5'-C4'	-2.04	1.45	1.51

All (44) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	418	5CA	O3'-C3'-C4'	8.83	136.44	111.08
4	A	418	5CA	O5'-C5'-C4'	-6.62	95.76	107.57
4	B	416	5CA	O3'-C3'-C4'	5.99	128.29	111.08
4	A	418	5CA	C5'-C4'-C3'	5.88	136.36	115.21
4	B	416	5CA	C5'-C4'-C3'	4.75	132.32	115.21
4	B	416	5CA	C4-C5-N7	-4.61	105.31	110.58
4	A	418	5CA	O2S-S-O1S	-4.29	114.43	120.85
4	B	416	5CA	O4'-C4'-C5'	4.26	122.99	109.33
4	B	416	5CA	C5-C4-N3	-4.25	120.86	126.72
4	B	416	5CA	O2S-S-O1S	-4.13	114.66	120.85
4	A	418	5CA	C2'-C3'-C4'	-4.06	94.76	102.61
4	A	418	5CA	N3-C2-N1	-4.06	122.44	128.58
4	A	418	5CA	CB-CA-C	3.98	118.14	109.52
4	B	416	5CA	N3-C2-N1	-3.95	122.60	128.58
4	B	416	5CA	O5'-C5'-C4'	-3.78	100.82	107.57
4	B	416	5CA	O4'-C4'-C3'	-3.72	97.77	105.15
4	A	418	5CA	C4'-O4'-C1'	-3.71	101.27	109.47
4	A	418	5CA	O3'-C3'-C2'	3.70	123.69	111.82
4	B	416	5CA	C2-N1-C6	3.60	124.65	118.73
4	A	418	5CA	O4'-C4'-C5'	3.57	120.76	109.33
4	B	416	5CA	O4'-C1'-N9	3.49	114.80	108.09
4	A	418	5CA	C2-N1-C6	3.43	124.37	118.73
4	A	418	5CA	C3'-C2'-C1'	-3.36	95.11	101.46
4	B	416	5CA	C3'-C2'-C1'	-3.23	95.36	101.46
4	B	416	5CA	C-CA-N	3.15	121.27	109.37
4	A	418	5CA	N6-C6-N1	3.02	125.10	118.38
4	A	418	5CA	O-C-CA	3.02	126.63	120.28
4	B	416	5CA	N3-C4-N9	2.93	132.14	127.17
4	A	418	5CA	C4-N9-C1'	-2.87	119.92	126.63
4	B	416	5CA	C5-N7-C8	2.77	107.81	103.45
6	B	417	EPE	O1S-S-C10	-2.77	102.54	106.73
6	B	417	EPE	C7-N4-C5	-2.75	103.91	111.24
4	A	418	5CA	C5-C4-N3	-2.74	122.94	126.72
4	B	416	5CA	O-C-N3S	2.74	127.96	122.98
4	A	418	5CA	C-CA-N	2.56	119.03	109.37
4	A	418	5CA	O-C-N3S	2.55	127.62	122.98
4	A	418	5CA	C5'-O5'-S	2.38	121.96	116.97

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	418	5CA	C1'-N9-C8	2.36	132.34	127.09
4	A	418	5CA	C4-C5-N7	-2.36	107.88	110.58
4	B	416	5CA	O-C-CA	2.33	125.18	120.28
4	B	416	5CA	CB-CA-C	2.27	114.42	109.52
4	A	418	5CA	C2'-C1'-N9	2.23	118.85	113.30
4	B	416	5CA	C6-C5-C4	2.22	120.21	117.18
4	A	418	5CA	N3-C4-N9	2.16	130.84	127.17

All (3) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
4	A	418	5CA	C4'
4	A	418	5CA	CA
4	B	416	5CA	CA

All (16) torsion outliers are listed below:

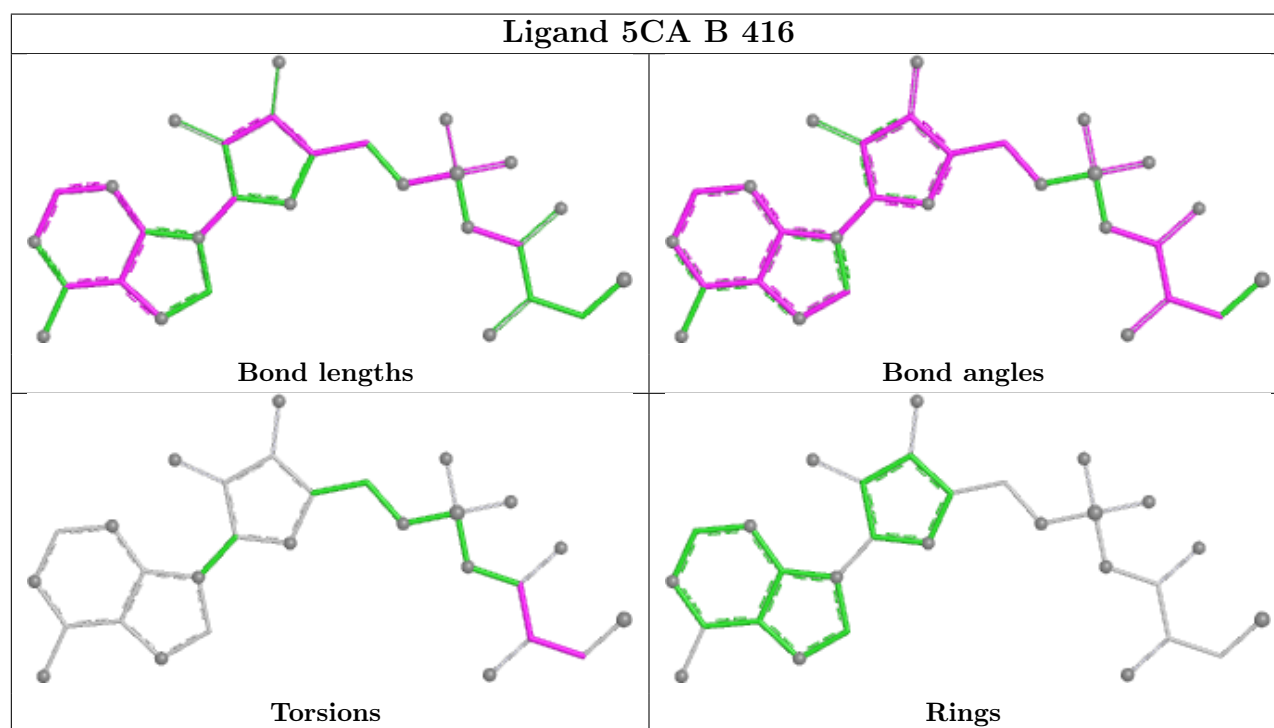
Mol	Chain	Res	Type	Atoms
4	A	418	5CA	N-CA-CB-SG
4	A	418	5CA	C3'-C4'-C5'-O5'
4	B	416	5CA	N-CA-CB-SG
6	B	417	EPE	C10-C9-N1-C6
5	A	419	1PE	OH6-C15-C25-OH5
6	B	417	EPE	N4-C7-C8-O8
5	A	419	1PE	OH2-C12-C22-OH3
5	A	419	1PE	OH7-C16-C26-OH6
4	A	418	5CA	O-C-CA-CB
5	A	419	1PE	C16-C26-OH6-C15
4	A	418	5CA	N3S-C-CA-N
4	A	418	5CA	N3S-C-CA-CB
4	B	416	5CA	O-C-CA-CB
4	B	416	5CA	N3S-C-CA-CB
5	A	419	1PE	C25-C15-OH6-C26
4	A	418	5CA	O4'-C4'-C5'-O5'

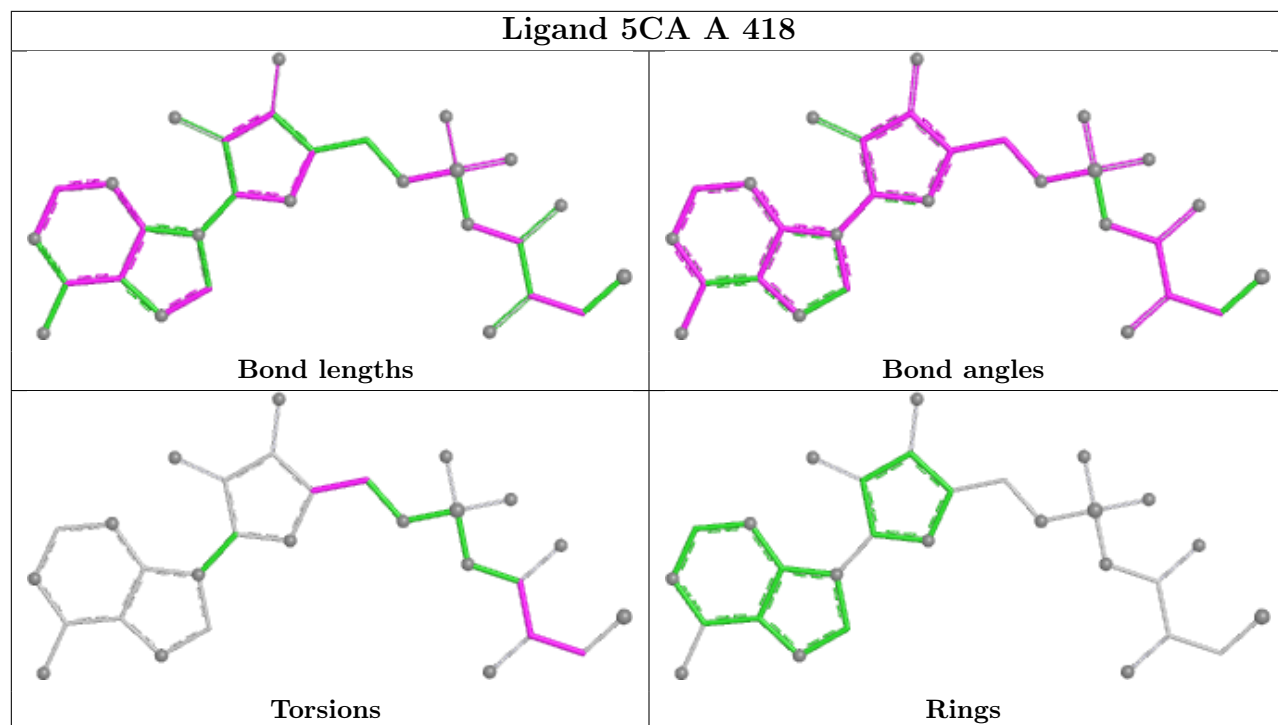
There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	419	1PE	1	0
4	B	416	5CA	1	0
4	A	418	5CA	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

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	399/414 (96%)	0.48	27 (6%) 23 24	14, 24, 36, 44	0
1	B	401/414 (96%)	0.12	24 (5%) 27 28	11, 18, 32, 48	0
All	All	800/828 (96%)	0.30	51 (6%) 25 26	11, 21, 35, 48	0

All (51) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	50	ALA	5.5
1	B	50	ALA	4.5
1	B	49	ASP	4.4
1	B	98	ILE	4.4
1	B	94	GLU	4.2
1	B	97	GLY	4.0
1	B	242	GLY	3.9
1	A	357	ASP	3.6
1	A	49	ASP	3.4
1	A	385	ASP	3.4
1	A	390	GLY	3.3
1	A	394	THR	3.3
1	A	327	PHE	3.3
1	B	392	HIS	3.3
1	B	327	PHE	3.2
1	B	331	GLU	3.2
1	B	99	ASP	3.1
1	B	96	ASP	3.1
1	B	93	ALA	3.0
1	A	298	VAL	2.8
1	A	37	THR	2.8
1	A	392	HIS	2.8
1	B	95	ARG	2.6
1	A	27	ARG	2.6

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	B	398	ARG	2.6
1	A	389	TYR	2.6
1	B	386	ALA	2.6
1	A	153	GLU	2.5
1	B	91	GLU	2.5
1	B	87	ASP	2.5
1	B	385	ASP	2.4
1	B	389	TYR	2.4
1	A	269	GLY	2.3
1	A	399	LEU	2.3
1	A	354	ASP	2.3
1	A	386	ALA	2.3
1	A	331	GLU	2.3
1	B	342	ARG	2.3
1	B	350	PRO	2.2
1	A	267	VAL	2.2
1	B	90	PHE	2.2
1	A	36	PRO	2.2
1	A	383	CYS	2.1
1	B	412	LEU	2.1
1	B	48	TYR	2.1
1	A	12	VAL	2.1
1	A	145	ALA	2.1
1	A	324	GLU	2.1
1	A	193	ASP	2.1
1	A	305	ALA	2.0
1	A	348	ALA	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,

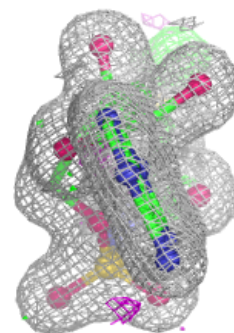
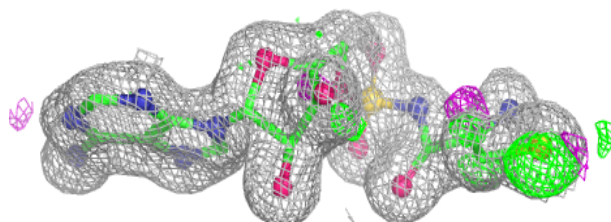
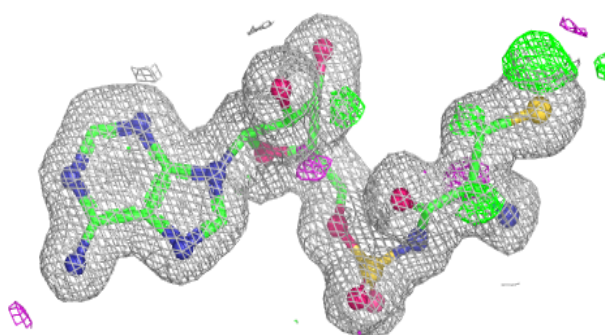
median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

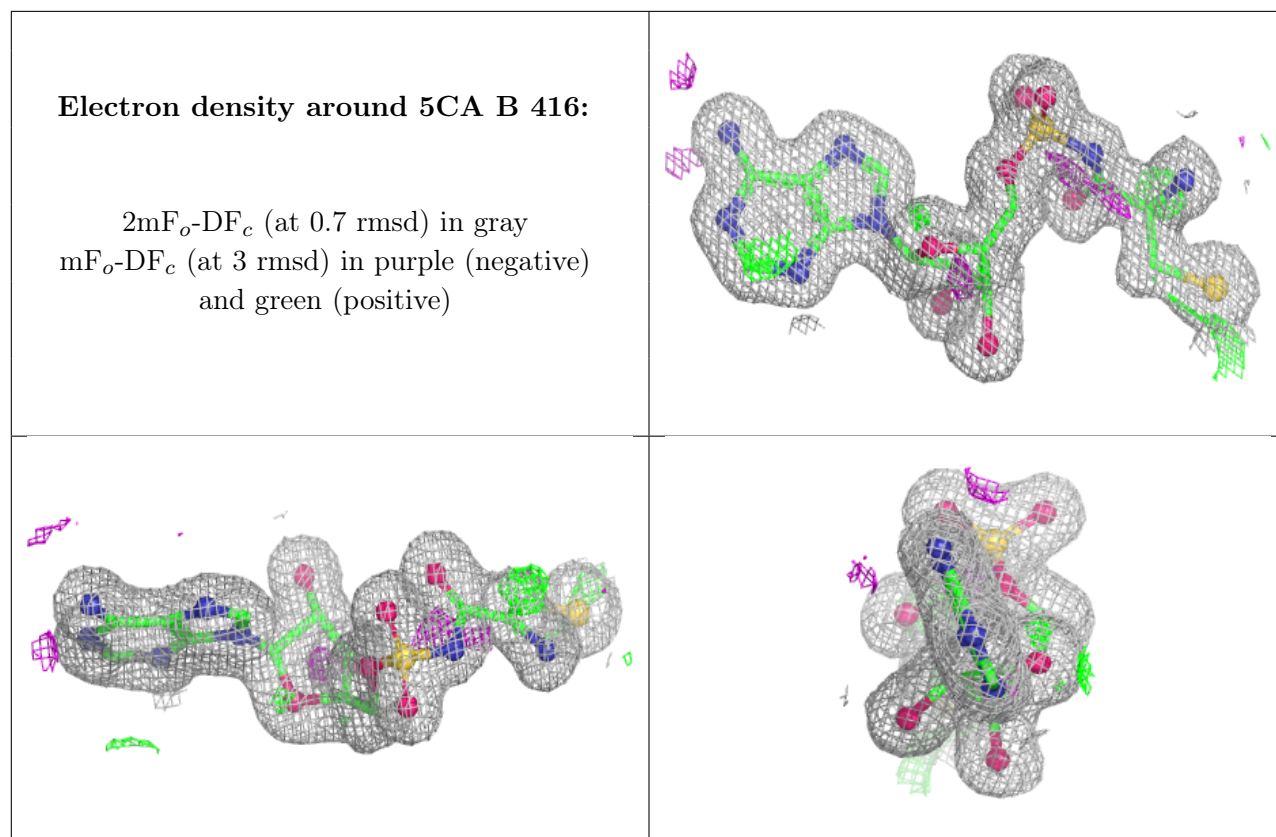
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
6	EPE	B	417	15/15	0.80	0.16	33,35,41,41	0
3	SO4	A	416	5/5	0.81	0.15	59,59,60,60	0
3	SO4	A	414	5/5	0.84	0.15	64,64,64,64	0
3	SO4	A	417	5/5	0.85	0.13	67,68,68,68	0
3	SO4	B	415	5/5	0.87	0.12	47,48,49,49	0
5	1PE	A	419	16/16	0.88	0.15	37,39,43,44	0
3	SO4	B	414	5/5	0.93	0.10	33,34,37,39	0
3	SO4	A	415	5/5	0.95	0.08	39,40,40,40	0
4	5CA	A	418	29/29	0.97	0.06	15,17,18,19	0
4	5CA	B	416	29/29	0.98	0.05	12,14,17,17	0
2	ZN	A	413	1/1	0.99	0.03	16,16,16,16	0
2	ZN	B	413	1/1	1.00	0.03	13,13,13,13	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.

Electron density around 5CA A 418:

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 ⓘ

There are no such residues in this entry.