



# Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 01:46 PM GMT

PDB ID : 1F4C  
Title : CRYSTAL STRUCTURE OF E. COLI THYMIDYLATE SYNTHASE CO-VALENTLY MODIFIED AT C146 WITH N-[TOSYL-D-PROLINYL]AMIN O-ETHANETHIOL  
Authors : Erlanson, D.A.; Braisted, A.C.; Raphael, D.R.; Randal, M.; Stroud, R.M.; Gordon, E.; Wells, J.A.  
Deposited on : 2000-06-07  
Resolution : 2.00 Å(reported)

This is a full wwPDB 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 <http://wwpdb.org/ValidationPDFNotes.html>

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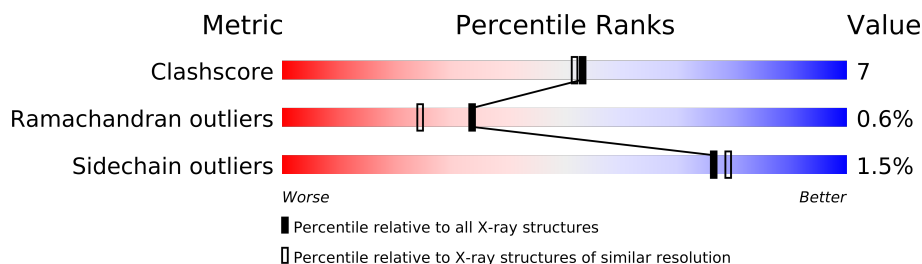
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : **NOT EXECUTED**  
EDS : **NOT EXECUTED**  
Percentile statistics : 21963  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance

The reported resolution of this entry is 2.00 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	79885	6188 (2.00-2.00)
Ramachandran outliers	78287	6102 (2.00-2.00)
Sidechain outliers	78261	6100 (2.00-2.00)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	264	
1	B	264	

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 4876 atoms, of which 0 are hydrogen and 0 are deuterium.

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.

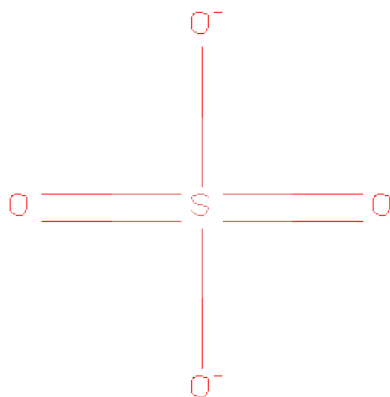
- Molecule 1 is a protein called THYMIDYLATE SYNTHASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	262	Total	C	N	O	S	0	2	0
			2149	1374	369	394	12			
1	B	261	Total	C	N	O	S	0	3	0
			2145	1370	369	394	12			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	CXM	MET	ENGINEERED	UNP P0A884
B	1	CXM	MET	ENGINEERED	UNP P0A884

- Molecule 2 is SULFATE ION (three-letter code: SO<sub>4</sub>) (formula: O<sub>4</sub>S).



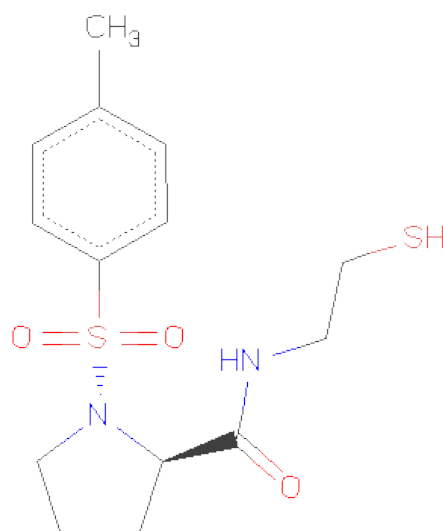
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		

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

- Molecule 3 is N-[TOSYL-D-PROLINYL]AMINO-ETHANETHIOL (three-letter code: TP2) (formula:  $C_{14}H_{20}N_2O_3S_2$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	S	0	0
			21	14	2	3	2		
3	B	1	Total	C	N	O	S	0	0
			21	14	2	3	2		

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula:  $C_3H_8O_3$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	C	O	0	0
			6	3	3		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	286	Total	O	0	0
			286	286		
5	B	223	Total	O	0	0
			223	223		

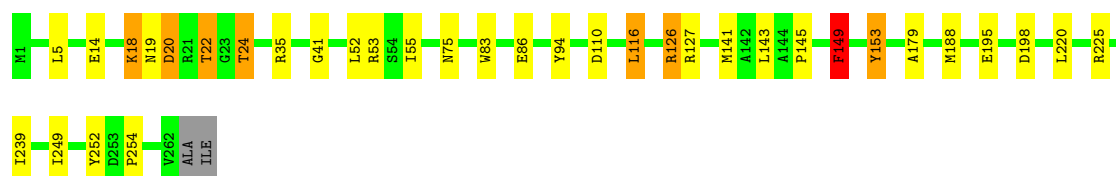
### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

Note EDS was not executed.

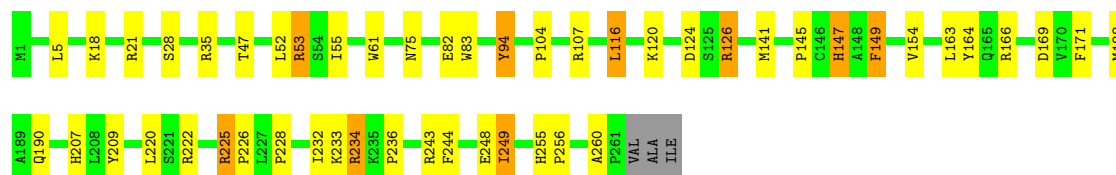
#### • Molecule 1: THYMIDYLATE SYNTHASE

Chain A: 



#### • Molecule 1: THYMIDYLATE SYNTHASE

Chain B: 



## 4 Data and refinement statistics

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

Property	Value	Source
Space group	P 63	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	126.22Å 126.22Å 67.02Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	10.00 – 2.00	Depositor
% Data completeness (in resolution range)	98.8 (10.00-2.00)	Depositor
$R_{merge}$	0.04	Depositor
$R_{sym}$	(Not available)	Depositor
Refinement program	REFMAC	Depositor
R, $R_{free}$	0.198 , 0.268	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	4876	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	34.0	wwPDB-VP

## 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: GOL, CXM, TP2, SO4

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	0.57	0/2209	1.34	18/3000 (0.6%)
1	B	0.51	0/2210	1.27	19/3001 (0.6%)
All	All	0.54	0/4419	1.31	37/6001 (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	2
1	B	0	1
All	All	0	3

There are no bond length outliers.

All (37) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	18	LYS	CA-CB-CG	13.58	143.28	113.40
1	A	127	ARG	NE-CZ-NH2	-12.68	113.96	120.30
1	A	127	ARG	NE-CZ-NH1	12.41	126.51	120.30
1	B	225	ARG	NE-CZ-NH1	10.75	125.68	120.30
1	B	35	ARG	NE-CZ-NH2	-8.81	115.90	120.30
1	B	107	ARG	CD-NE-CZ	8.37	135.32	123.60
1	B	243	ARG	NE-CZ-NH1	8.34	124.47	120.30
1	B	222	ARG	NE-CZ-NH2	-8.15	116.22	120.30
1	A	225	ARG	NE-CZ-NH1	7.97	124.29	120.30
1	A	110	ASP	CB-CG-OD2	7.51	125.06	118.30
1	B	126	ARG	NE-CZ-NH2	-7.33	116.63	120.30
1	A	225	ARG	CD-NE-CZ	7.18	133.65	123.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	225	ARG	NE-CZ-NH2	-7.06	116.77	120.30
1	B	61	TRP	CA-CB-CG	7.03	127.05	113.70
1	A	153	TYR	CB-CG-CD2	-6.77	116.94	121.00
1	B	234	ARG	NE-CZ-NH2	6.49	123.54	120.30
1	A	126	ARG	NE-CZ-NH1	-6.46	117.07	120.30
1	B	107	ARG	NE-CZ-NH1	6.40	123.50	120.30
1	B	234	ARG	NE-CZ-NH1	-6.37	117.11	120.30
1	B	149[A]	PHE	CA-CB-CG	6.14	128.64	113.90
1	B	149[B]	PHE	CA-CB-CG	6.14	128.64	113.90
1	B	107	ARG	NE-CZ-NH2	-6.14	117.23	120.30
1	A	225	ARG	CG-CD-NE	6.08	124.58	111.80
1	A	86	GLU	OE1-CD-OE2	-5.87	116.26	123.30
1	B	53	ARG	NE-CZ-NH2	-5.85	117.37	120.30
1	A	195	GLU	OE1-CD-OE2	-5.79	116.34	123.30
1	A	153	TYR	CB-CG-CD1	5.63	124.38	121.00
1	A	149[A]	PHE	CA-CB-CG	5.57	127.26	113.90
1	A	149[B]	PHE	CA-CB-CG	5.57	127.26	113.90
1	A	126	ARG	NE-CZ-NH2	-5.51	117.55	120.30
1	A	126	ARG	NH1-CZ-NH2	5.44	125.38	119.40
1	B	169	ASP	CB-CG-OD1	5.40	123.16	118.30
1	B	21	ARG	NE-CZ-NH1	5.39	123.00	120.30
1	B	94	TYR	CB-CG-CD2	-5.17	117.90	121.00
1	B	94	TYR	CB-CG-CD1	5.08	124.05	121.00
1	A	18	LYS	CB-CA-C	5.06	120.53	110.40
1	B	243	ARG	NE-CZ-NH2	-5.04	117.78	120.30

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	14	GLU	Mainchain
1	A	41	GLY	Mainchain
1	B	147	HIS	Mainchain

## 5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2149	0	2070	30	0
1	B	2145	0	2063	44	0
2	A	15	0	0	1	0
2	B	10	0	0	0	0
3	A	21	0	19	0	0
3	B	21	0	19	1	0
4	B	6	0	8	1	0
5	A	286	0	0	3	0
5	B	223	0	0	2	0
All	All	4876	0	4179	60	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 7.

All (60) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:149[A]:PHE:CE2	1:B:164:TYR:CB	2.55	0.89
1:A:149[A]:PHE:HZ	1:B:164:TYR:CB	1.87	0.86
1:A:149[A]:PHE:HZ	1:B:164:TYR:CG	1.94	0.86
1:B:166:ARG:HH12	4:B:702:GOL:H2	1.43	0.83
1:A:149[A]:PHE:CZ	1:B:164:TYR:CG	2.71	0.78
1:B:149[A]:PHE:CE2	1:B:164:TYR:HB3	2.23	0.73
1:B:149[A]:PHE:CE2	1:B:164:TYR:HB2	2.24	0.71
1:A:22:THR:HA	5:A:1080:HOH:O	1.91	0.70
1:B:149[A]:PHE:CD2	1:B:164:TYR:HB3	2.27	0.68
1:A:18:LYS:HB3	5:A:900:HOH:O	1.96	0.65
1:A:20:ASP:HB2	1:B:126:ARG:HH12	1.63	0.63
1:A:149[A]:PHE:HZ	1:B:164:TYR:HB3	1.65	0.61
1:B:190:GLN:HE22	1:B:232:ILE:HG21	1.65	0.61
1:A:52:LEU:HD22	1:A:249:ILE:HG13	1.85	0.59
1:B:147:HIS:HB2	1:B:163:LEU:HD11	1.85	0.58
1:B:83:TRP:CZ3	3:B:802:TP2:H202	2.38	0.58
1:B:141:MET:SD	1:B:145:PRO:HD3	2.44	0.57
1:A:83:TRP:CZ3	1:A:143:LEU:HD21	2.40	0.56
1:A:149[A]:PHE:CE2	1:B:164:TYR:CD2	2.94	0.56
1:B:82:GLU:HG2	5:B:1007:HOH:O	2.04	0.56
1:A:18:LYS:HE3	1:B:154:VAL:O	2.06	0.56
1:A:20:ASP:OD1	1:A:24:THR:HB	2.05	0.55
1:A:19:ASN:HB2	5:B:1010:HOH:O	2.07	0.53
1:B:255:HIS:HB3	1:B:256:PRO:HD2	1.90	0.52
1:B:149[A]:PHE:CD2	1:B:164:TYR:CB	2.92	0.52
1:B:28:SER:HB3	1:B:207:HIS:HB3	1.92	0.51
1:B:116:LEU:HD23	1:B:120:LYS:HE2	1.94	0.49

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:18:LYS:HE2	1:B:124:ASP:OD1	2.13	0.49
1:A:20:ASP:HB2	1:B:126:ARG:NH1	2.28	0.49
1:B:52:LEU:HA	1:B:55:ILE:HD12	1.93	0.48
1:A:35:ARG:HD3	1:A:198:ASP:OD2	2.13	0.48
1:A:5:LEU:HD22	1:A:220:LEU:HD23	1.96	0.48
1:A:126:ARG:NH2	1:B:209:TYR:OH	2.46	0.48
1:A:149[A]:PHE:HE2	1:B:164:TYR:CD2	2.32	0.47
1:A:252:TYR:CE2	1:A:254:PRO:HG3	2.49	0.47
1:A:116:LEU:HD23	1:A:239:ILE:HG21	1.96	0.47
1:B:225:ARG:HB3	1:B:226:PRO:HD2	1.95	0.47
1:B:234:ARG:O	1:B:236:PRO:HD3	2.14	0.47
1:A:149[A]:PHE:CZ	1:B:164:TYR:CD2	3.02	0.47
1:B:82:GLU:HG3	1:B:83:TRP:CE3	2.50	0.46
1:B:53:ARG:HG3	1:B:244:PHE:CE1	2.51	0.46
1:A:141:MET:SD	1:A:145:PRO:HD3	2.56	0.46
1:A:53:ARG:NH1	1:A:75:ASN:O	2.39	0.46
1:B:5:LEU:HD22	1:B:220:LEU:HD23	1.98	0.45
1:B:171:PHE:CE2	1:B:260:ALA:HB2	2.52	0.45
1:B:18:LYS:HD2	1:B:28:SER:OG	2.16	0.44
1:B:233:LYS:HD2	1:B:248:GLU:CG	2.47	0.44
1:A:153:TYR:CZ	1:B:18:LYS:HE3	2.53	0.44
1:A:55:ILE:HD13	1:A:179:ALA:HB3	1.99	0.44
1:B:233:LYS:HD2	1:B:248:GLU:HG3	2.01	0.42
1:A:53:ARG:HB2	2:A:711:SO4:O2	2.19	0.42
1:B:255:HIS:HB3	1:B:256:PRO:CD	2.49	0.41
1:A:20:ASP:HB2	1:B:126:ARG:HH22	1.86	0.41
1:A:116:LEU:HD22	1:A:188:MET:CE	2.51	0.41
1:A:116:LEU:HD22	1:A:188:MET:HE1	2.02	0.41
1:B:5:LEU:HD11	1:B:47:THR:HG21	2.03	0.41
1:B:228:PRO:HB2	1:B:249:ILE:HD12	2.03	0.41
5:A:1003:HOH:O	1:B:104:PRO:HG2	2.21	0.41
1:B:225:ARG:HB3	1:B:226:PRO:CD	2.50	0.40
1:B:53:ARG:NH1	1:B:75:ASN:O	2.54	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	262/264 (99%)	252 (96%)	8 (3%)	2 (1%)	27	17
1	B	262/264 (99%)	254 (97%)	7 (3%)	1 (0%)	43	36
All	All	524/528 (99%)	506 (97%)	15 (3%)	3 (1%)	33	24

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	94	TYR
1	B	94	TYR
1	A	24	THR

### 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	233/232 (100%)	228 (98%)	5 (2%)	66	67
1	B	233/232 (100%)	230 (99%)	3 (1%)	80	82
All	All	466/464 (100%)	458 (98%)	8 (2%)	76	75

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	20	ASP
1	A	22	THR
1	A	116	LEU
1	A	149[A]	PHE
1	A	149[B]	PHE
1	B	116	LEU
1	B	188	MET
1	B	249	ILE

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

Mol	Chain	Res	Type
1	A	117	ASN
1	A	219	GLN
1	B	17	GLN
1	B	190	GLN
1	B	219	GLN

### 5.3.3 RNA ⓘ

There are no RNA chains in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

2 non-standard protein/DNA/RNA residues 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$
1	CXM	A	1	1	10,10,11	7.68	5 (50%)	8,11,13	8.08	3 (37%)
1	CXM	B	1	1	10,10,11	7.20	3 (30%)	8,11,13	18.41	5 (62%)

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
1	CXM	A	1	1	-	0/8/10/12	0/0/0/0
1	CXM	B	1	1	-	0/8/10/12	0/0/0/0

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	1	CXM	O-C	22.65	1.27	1.11
1	B	1	CXM	O-C	22.09	1.26	1.11
1	A	1	CXM	CA-N	-5.66	1.38	1.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	1	CXM	CB-CA	4.12	1.56	1.53
1	B	1	CXM	ON1-CN	3.72	1.28	1.21
1	A	1	CXM	CN-N	3.47	1.37	1.33
1	A	1	CXM	CA-C	3.45	1.54	1.48
1	B	1	CXM	CA-C	2.90	1.53	1.48

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	1	CXM	ON2-CN-ON1	-51.65	57.71	122.17
1	A	1	CXM	ON2-CN-ON1	-17.89	99.84	122.17
1	A	1	CXM	ON2-CN-N	-10.47	103.30	116.33
1	A	1	CXM	CA-N-CN	-9.27	107.50	121.67
1	B	1	CXM	ON1-CN-N	-4.20	117.41	124.90
1	B	1	CXM	CB-CA-N	3.73	115.64	110.48
1	B	1	CXM	CA-N-CN	-2.57	117.74	121.67
1	B	1	CXM	ON2-CN-N	-2.35	113.41	116.33

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

## 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

8 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	$\# Z  > 2$
2	SO4	A	701	-	4,4,4	0.80	0	6,6,6	0.29	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	SO4	A	711	-	4,4,4	1.00	0	6,6,6	0.35	0
2	SO4	A	712	-	4,4,4	0.95	0	6,6,6	0.57	0
3	TP2	A	801	1	22,22,22	4.12	3 (13%)	31,31,31	1.65	6 (19%)
4	GOL	B	702	-	5,5,5	0.57	0	5,5,5	0.46	0
2	SO4	B	713	-	4,4,4	0.99	0	6,6,6	0.49	0
2	SO4	B	714	-	4,4,4	0.93	0	6,6,6	0.22	0
3	TP2	B	802	1	22,22,22	3.88	3 (13%)	31,31,31	1.53	7 (22%)

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
2	SO4	A	701	-	-	0/0/0/0	0/0/0/0
2	SO4	A	711	-	-	0/0/0/0	0/0/0/0
2	SO4	A	712	-	-	0/0/0/0	0/0/0/0
3	TP2	A	801	1	-	0/20/30/30	0/2/2/2
4	GOL	B	702	-	-	0/4/4/4	0/0/0/0
2	SO4	B	713	-	-	0/0/0/0	0/0/0/0
2	SO4	B	714	-	-	0/0/0/0	0/0/0/0
3	TP2	B	802	1	-	0/20/30/30	0/2/2/2

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	801	TP2	O16-S8	12.84	1.59	1.43
3	A	801	TP2	O17-S8	12.12	1.59	1.43
3	B	802	TP2	O16-S8	11.97	1.58	1.43
3	B	802	TP2	O17-S8	11.06	1.57	1.43
3	B	802	TP2	S8-N7	7.56	1.74	1.63
3	A	801	TP2	S8-N7	7.05	1.74	1.63

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	801	TP2	O17-S8-O16	-5.52	109.53	119.38
3	A	801	TP2	C9-S8-N7	3.66	113.57	107.44
3	B	802	TP2	O17-S8-O16	-3.29	113.52	119.38
3	B	802	TP2	C3-C2-S1	-3.20	105.10	113.50
3	B	802	TP2	C6-N7-S8	-3.00	114.79	119.67
3	A	801	TP2	C3-C2-S1	-2.83	106.07	113.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	802	TP2	C9-S8-N7	2.73	112.01	107.44
3	A	801	TP2	C6-C5-N4	2.65	121.59	116.12
3	A	801	TP2	C6-N7-S8	-2.59	115.46	119.67
3	A	801	TP2	O21-C5-C6	-2.37	115.12	120.38
3	B	802	TP2	C14-C9-C10	2.28	124.00	120.39
3	B	802	TP2	C13-C14-C9	-2.09	117.10	119.50
3	B	802	TP2	C20-C6-C5	2.07	114.87	111.15

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

## 5.7 Other polymers ⓘ

There are no such residues in this entry.

## 5.8 Polymer linkage issues

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 will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

### 6.3 Carbohydrates ⓘ

EDS was not executed - this section will therefore be empty.

### 6.4 Ligands ⓘ

EDS was not executed - this section will therefore be empty.

### 6.5 Other polymers ⓘ

EDS was not executed - this section will therefore be empty.