



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 03:17 AM UTC

PDB ID : 1F4C / pdb_00001f4c
Title : CRYSTAL STRUCTURE OF E. COLI THYMIDYLATE SYNTHASE CO-VALENTLY MODIFIED AT C146 WITH N-[TOSYL-D-PROLINYL]AMIN O-ETHANETHIOL
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Deposited on : 2000-06-07
Resolution : 2.00 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

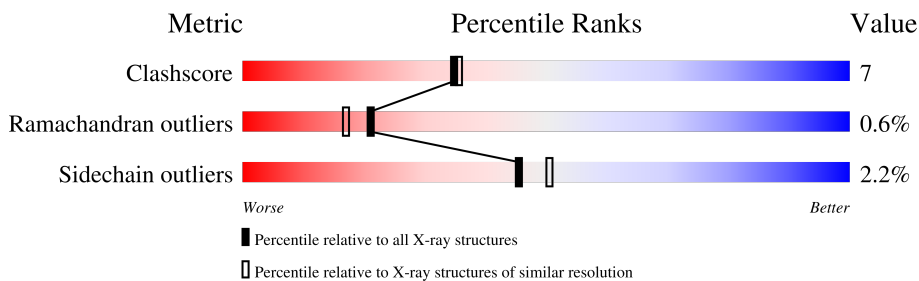
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 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	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$.

Note EDS was not executed.

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

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 4876 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 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 mutation	UNP P0A884
B	1	CXM	MET	engineered mutation	UNP P0A884

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



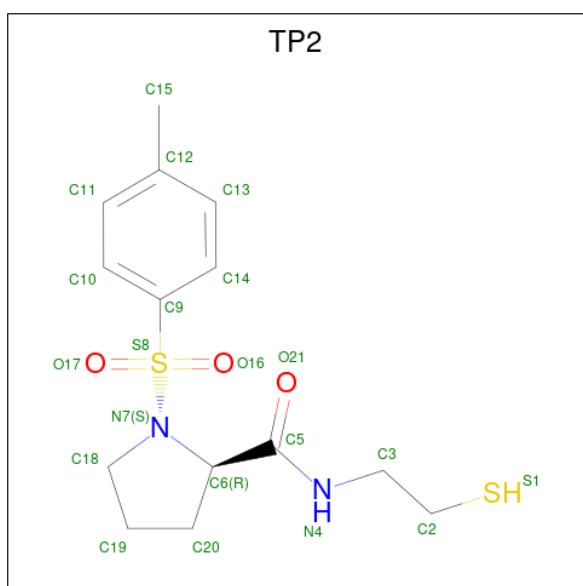
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	O S	0	0
			5	4 1		

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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 (CCD ID: 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 (CCD ID: 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		

Note EDS was not executed.

Chain A: 82% 14% ...



Chain B: 76% 21% ..



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 63	Depositor
Cell constants a, b, c, α , β , γ	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 ²)	34.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: TP2, SO4, GOL, CXM

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	0.76	0/2209	1.62	24/3000 (0.8%)
1	B	0.65	0/2210	1.59	17/3001 (0.6%)
All	All	0.70	0/4419	1.60	41/6001 (0.7%)

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 (41) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	18	LYS	CA-CB-CG	14.59	143.28	114.10
1	B	149[A]	PHE	CA-CB-CG	14.28	128.08	113.80
1	B	149[B]	PHE	CA-CB-CG	14.28	128.08	113.80
1	B	62	PHE	CA-CB-CG	9.44	123.24	113.80
1	B	150	PHE	CA-CB-CG	9.19	122.99	113.80
1	A	149[A]	PHE	CA-CB-CG	9.01	122.81	113.80
1	A	149[B]	PHE	CA-CB-CG	9.01	122.81	113.80
1	B	107	ARG	CD-NE-CZ	7.80	135.32	124.40
1	B	61	TRP	CA-CB-CG	7.08	127.05	113.60
1	A	62	PHE	CA-CB-CG	6.82	120.62	113.80
1	A	225	ARG	CD-NE-CZ	6.61	133.65	124.40
1	A	147	HIS	CA-CB-CG	6.48	120.28	113.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	73	HIS	CA-CB-CG	-6.48	107.32	113.80
1	A	20	ASP	CA-CB-CG	6.25	118.85	112.60
1	B	66	ASP	CA-CB-CG	6.25	118.85	112.60
1	A	19	ASN	CA-CB-CG	6.05	118.65	112.60
1	A	198	ASP	CA-CB-CG	6.03	118.63	112.60
1	A	143	LEU	CA-C-O	-6.03	113.12	120.54
1	A	211	ASN	CA-CB-CG	5.97	118.57	112.60
1	A	127	ARG	NE-CZ-NH2	-5.82	113.96	119.20
1	B	213	MET	N-CA-C	5.80	117.69	111.36
1	A	134	ASN	CA-CB-CG	-5.78	106.83	112.60
1	B	198	ASP	CA-CB-CG	5.74	118.34	112.60
1	A	225	ARG	CG-CD-NE	5.72	124.58	112.00
1	A	170	VAL	N-CA-C	5.70	115.89	110.42
1	A	174	LEU	N-CA-C	5.70	120.65	113.25
1	B	174	LEU	N-CA-C	5.50	120.40	113.25
1	A	213	MET	N-CA-C	5.44	116.89	111.07
1	B	146	CYS	CA-C-O	-5.38	113.43	120.00
1	B	19	ASN	CA-C-O	-5.35	115.13	121.66
1	A	207	HIS	N-CA-C	5.33	117.78	109.52
1	B	217	HIS	CA-CB-CG	-5.30	108.50	113.80
1	A	14	GLU	CB-CA-C	-5.13	100.46	109.07
1	A	149[A]	PHE	CB-CA-C	5.10	119.50	110.16
1	A	149[B]	PHE	CB-CA-C	5.10	119.50	110.16
1	B	255	HIS	CA-CB-CG	5.07	118.87	113.80
1	A	122	ASP	CA-C-N	5.03	124.69	119.56
1	A	122	ASP	C-N-CA	5.03	124.69	119.56
1	A	127	ARG	NE-CZ-NH1	5.01	126.51	121.50
1	B	202	THR	CA-C-N	-5.00	117.76	122.66
1	B	202	THR	C-N-CA	-5.00	117.76	122.66

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 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	2149	0	2070	31	0
1	B	2145	0	2063	47	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	64	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 7.

All (64) 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: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:83:TRP:CZ3	3:B:802:TP2:H202	2.38	0.58
1:B:147:HIS:HB2	1:B:163:LEU:HD11	1.85	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:116:LEU:HD22	1:A:188:MET:HE3	1.89	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
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:23:GLY:O	1:A:24:THR:C	2.62	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: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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:8:MET:HE1	1:B:219:GLN:NE2	2.36	0.41
1:B:124:ASP:O	1:B:125:SER:C	2.63	0.41
1:B:53:ARG:NH1	1:B:75:ASN:O	2.54	0.40
1:B:93:VAL:O	1:B:94:TYR:C	2.65	0.40
1:B:225:ARG:HB3	1:B:226:PRO:CD	2.50	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	262/264 (99%)	252 (96%)	8 (3%)	2 (1%)	16	11
1	B	262/264 (99%)	254 (97%)	7 (3%)	1 (0%)	30	27
All	All	524/528 (99%)	506 (97%)	15 (3%)	3 (1%)	21	17

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 [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	233/232 (100%)	225 (97%)	8 (3%)	32	33
1	B	233/232 (100%)	230 (99%)	3 (1%)	61	68
All	All	466/464 (100%)	455 (98%)	11 (2%)	45	47

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

Mol	Chain	Res	Type
1	A	20	ASP
1	A	22	THR
1	A	103	THR
1	A	115	VAL
1	A	116	LEU
1	A	149[A]	PHE
1	A	149[B]	PHE
1	A	237	GLU
1	B	116	LEU
1	B	188	MET
1	B	249	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	19	ASN
1	A	117	ASN
1	A	211	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 molecules 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	9,10,11	2.15	1 (11%)	9,11,13	2.94	2 (22%)
1	CXM	B	1	1	9,10,11	1.56	1 (11%)	9,11,13	3.67	6 (66%)

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	-	1/9/10/12	-
1	CXM	B	1	1	-	2/9/10/12	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	1	CXM	CA-N	-5.38	1.38	1.46
1	B	1	CXM	ON1-CN	3.76	1.28	1.21

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	1	CXM	ON2-CN-ON1	-8.36	57.71	121.85
1	A	1	CXM	CA-N-CN	-7.96	107.50	122.11
1	B	1	CXM	ON1-CN-N	-4.54	117.41	124.86
1	B	1	CXM	C-CA-N	-3.53	102.68	109.50
1	A	1	CXM	ON2-CN-ON1	-2.87	99.84	121.85
1	B	1	CXM	CB-CA-N	2.81	115.64	110.52
1	B	1	CXM	CA-N-CN	-2.38	117.74	122.11
1	B	1	CXM	O-C-CA	-2.02	119.56	124.77

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	B	1	CXM	ON2-CN-N-CA
1	B	1	CXM	O-C-CA-CB

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Mol	Chain	Res	Type	Atoms
1	A	1	CXM	ON1-CN-N-CA

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

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	711	-	4,4,4	0.72	0	6,6,6	0.36	0
4	GOL	B	702	-	5,5,5	0.71	0	5,5,5	0.55	0
3	TP2	A	801	1	21,22,22	4.92	3 (14%)	30,31,31	1.66	4 (13%)
2	SO4	A	712	-	4,4,4	0.69	0	6,6,6	0.34	0
2	SO4	B	714	-	4,4,4	0.66	0	6,6,6	0.33	0
2	SO4	B	713	-	4,4,4	0.61	0	6,6,6	0.29	0
3	TP2	B	802	1	21,22,22	4.61	3 (14%)	30,31,31	1.50	6 (20%)
2	SO4	A	701	-	4,4,4	0.53	0	6,6,6	0.34	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
4	GOL	B	702	-	-	2/4/4/4	-
3	TP2	A	801	1	-	0/20/30/30	0/2/2/2

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
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	15.34	1.59	1.43
3	A	801	TP2	O17-S8	14.49	1.59	1.43
3	B	802	TP2	O16-S8	14.31	1.58	1.43
3	B	802	TP2	O17-S8	13.23	1.57	1.43
3	B	802	TP2	S8-N7	7.70	1.74	1.63
3	A	801	TP2	S8-N7	7.18	1.74	1.63

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	801	TP2	O17-S8-O16	-6.45	109.53	119.59
3	B	802	TP2	O17-S8-O16	-3.89	113.52	119.59
3	A	801	TP2	C9-S8-N7	3.56	113.57	107.36
3	B	802	TP2	C14-C9-C10	2.71	124.00	120.47
3	B	802	TP2	C9-S8-N7	2.67	112.01	107.36
3	A	801	TP2	C6-C5-N4	2.60	121.59	115.99
3	B	802	TP2	C13-C14-C9	-2.40	117.10	119.44
3	A	801	TP2	O21-C5-C6	-2.31	115.12	120.69
3	B	802	TP2	C6-N7-S8	-2.20	114.79	119.77
3	B	802	TP2	C3-N4-C5	-2.05	118.87	122.55

There are no chirality outliers.

All (2) torsion outliers are listed below:

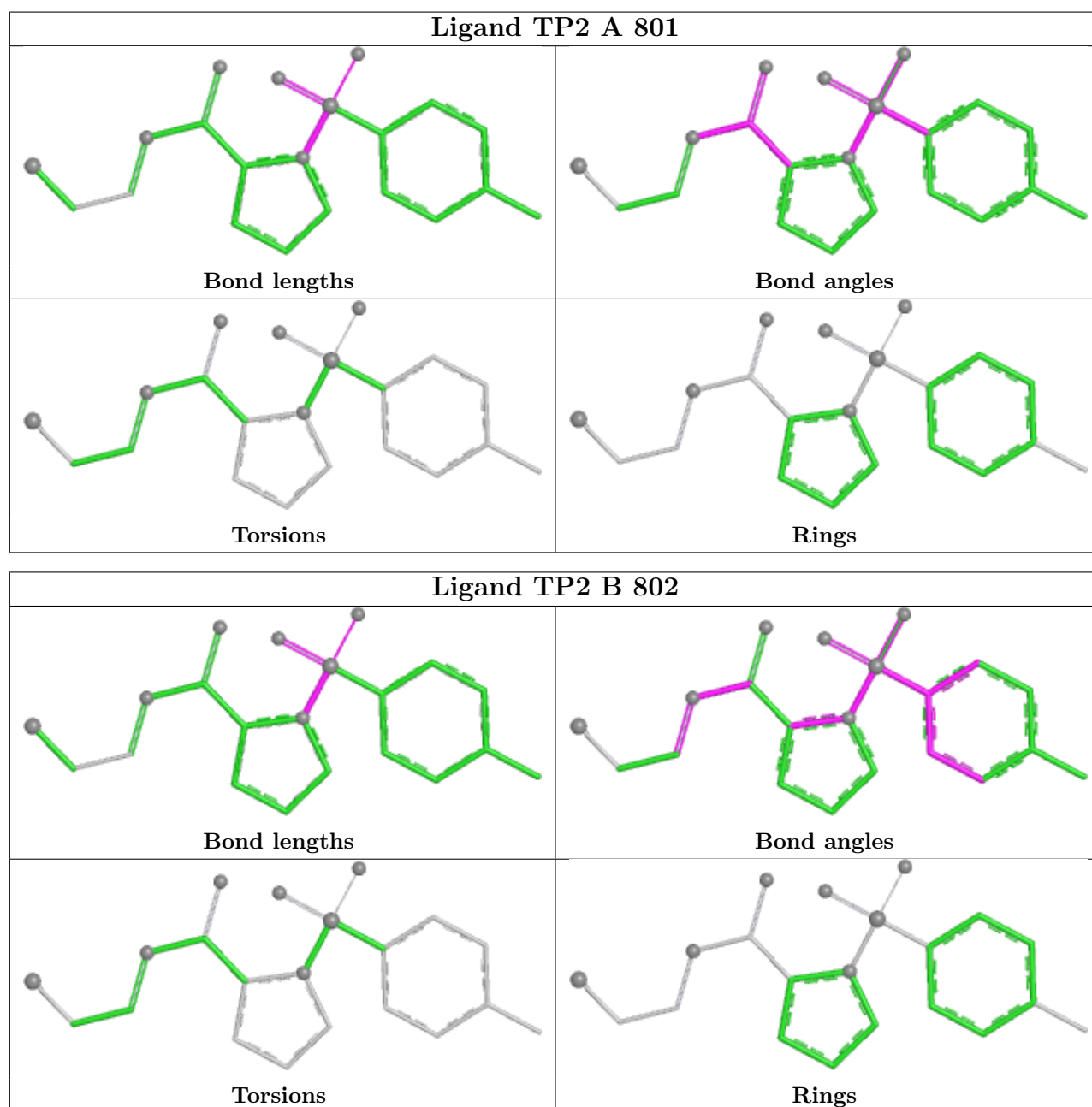
Mol	Chain	Res	Type	Atoms
4	B	702	GOL	C1-C2-C3-O3
4	B	702	GOL	O2-C2-C3-O3

There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	711	SO4	1	0
4	B	702	GOL	1	0
3	B	802	TP2	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

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.