



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 06:56 PM UTC

PDB ID : 1F4F / pdb_00001f4f
Title : CRYSTAL STRUCTURE OF E. COLI THYMIDYLATE SYNTHASE COM-
PLEXED WITH SP-722
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 X-ray Structure Validation Report for a publicly released PDB entry.

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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

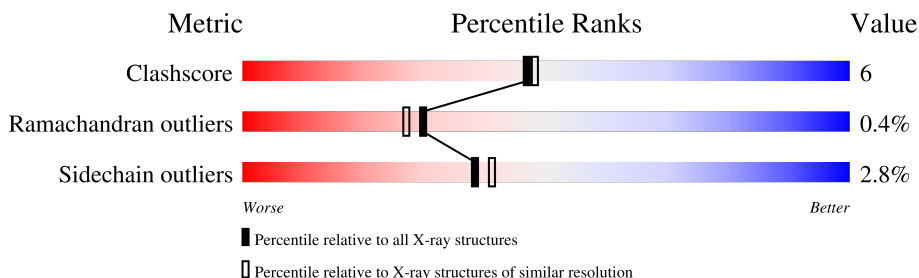
1 Overall quality at a glance

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	 81% 16% ...
1	B	264	 81% 17% ..

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 4842 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.

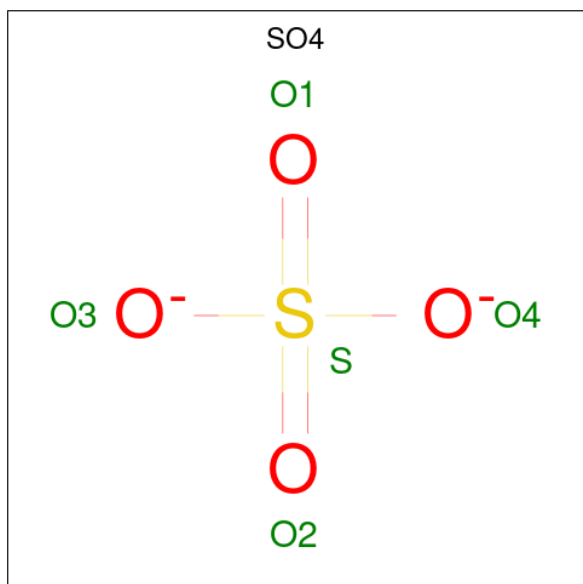
- Molecule 1 is a protein called THYMIDYLATE SYNTHASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	262	Total	C	N	O	S	0	1	0
			2143	1368	369	394	12			
1	B	261	Total	C	N	O	S	0	2	0
			2139	1364	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).



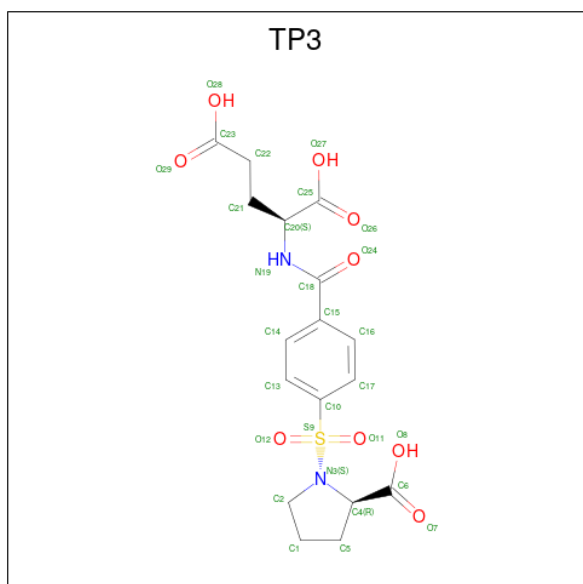
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		

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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		

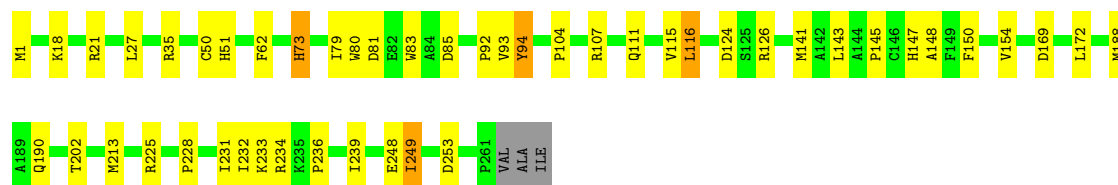
- Molecule 3 is 4-[[GLUTAMIC ACID]-CARBONYL]-BENZENE-SULFONYL-D-PROLINE (CCD ID: TP3) (formula: C₁₇H₂₀N₂O₉S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	S	0	0
			29	17	2	9	1		
3	B	1	Total	C	N	O	S	0	1
			33	19	2	11	1		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	266	Total	O	0	0
			266	266		
4	B	207	Total	O	1	0
			207	207		



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	126.14Å 126.14Å 66.81Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	10.00 – 2.00	Depositor
% Data completeness (in resolution range)	99.4 (10.00-2.00)	Depositor
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	REFMAC	Depositor
R, R_{free}	0.194 , 0.251	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	4842	wwPDB-VP
Average B, all atoms (Å ²)	37.0	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: CXM, TP3, 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	0.53	0/2197	1.38	14/2984 (0.5%)
1	B	0.49	0/2198	1.37	20/2985 (0.7%)
All	All	0.51	0/4395	1.38	34/5969 (0.6%)

There are no bond length outliers.

All (34) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	150	PHE	CA-CB-CG	8.68	122.48	113.80
1	A	225	ARG	CD-NE-CZ	8.28	135.99	124.40
1	B	150	PHE	CA-CB-CG	8.12	121.92	113.80
1	B	73	HIS	CA-CB-CG	-7.65	106.15	113.80
1	A	18	LYS	CA-CB-CG	7.27	128.65	114.10
1	A	62	PHE	CA-CB-CG	6.88	120.68	113.80
1	B	62	PHE	CA-CB-CG	6.70	120.50	113.80
1	A	176	PHE	CA-CB-CG	6.59	120.39	113.80
1	B	81	ASP	N-CA-C	6.28	118.93	111.33
1	A	148	ALA	N-CA-C	6.22	121.05	112.90
1	B	35	ARG	CD-NE-CZ	6.11	132.95	124.40
1	B	148	ALA	N-CA-C	6.01	120.89	113.50
1	A	225	ARG	CG-CD-NE	5.83	124.83	112.00
1	A	94	TYR	CA-C-N	5.77	126.49	120.03
1	A	94	TYR	C-N-CA	5.77	126.49	120.03
1	A	129	ILE	N-CA-C	5.75	118.23	108.86
1	B	147	HIS	CA-CB-CG	5.71	119.52	113.80
1	B	213	MET	N-CA-C	5.53	117.39	111.36
1	B	202	THR	CA-C-N	-5.48	117.29	122.66
1	B	202	THR	C-N-CA	-5.48	117.29	122.66
1	A	28	SER	N-CA-C	5.39	117.58	109.23
1	B	169	ASP	CA-CB-CG	5.39	117.99	112.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	50	CYS	CA-C-O	-5.28	115.19	121.16
1	B	27	LEU	CA-C-N	-5.25	114.35	122.59
1	B	27	LEU	C-N-CA	-5.25	114.35	122.59
1	A	20	ASP	CA-CB-CG	5.22	117.82	112.60
1	B	27	LEU	CA-C-O	-5.18	114.62	120.32
1	B	27	LEU	N-CA-C	-5.17	100.09	108.52
1	A	66	ASP	CA-CB-CG	5.14	117.74	112.60
1	B	107	ARG	CD-NE-CZ	5.10	131.54	124.40
1	B	85	ASP	CA-CB-CG	5.06	117.66	112.60
1	B	50	CYS	CA-C-N	-5.06	116.26	122.84
1	B	50	CYS	C-N-CA	-5.06	116.26	122.84
1	A	207	HIS	N-CA-C	5.04	116.85	109.24

There are no chirality outliers.

There are no planarity outliers.

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	2143	0	2066	30	0
1	B	2139	0	2059	28	0
2	A	15	0	0	0	0
2	B	10	0	0	0	0
3	A	29	0	17	1	0
3	B	33	0	8	1	0
4	A	266	0	0	3	0
4	B	207	0	0	2	0
All	All	4842	0	4150	49	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 6.

All (49) 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:20:ASP:HB2	1:B:126:ARG:HH12	1.41	0.83
1:A:116:LEU:HD22	1:A:188:MET:HE3	1.67	0.75
1:A:83:TRP:CZ3	1:A:143:LEU:HD21	2.33	0.63
1:A:52:LEU:HD22	1:A:249:ILE:HG13	1.82	0.61
1:A:79:ILE:HG22	3:A:701:TP3:H222	1.83	0.58
1:B:79:ILE:HG22	3:B:702[B]:TP3:H222	1.86	0.57
1:B:141:MET:SD	1:B:145:PRO:HD3	2.48	0.54
1:A:20:ASP:HB2	1:B:126:ARG:NH1	2.18	0.53
1:A:125:SER:HB2	1:B:21:ARG:HH21	1.74	0.53
1:B:83:TRP:CE2	1:B:143:LEU:HD11	2.44	0.53
1:A:141:MET:SD	1:A:145:PRO:HD3	2.50	0.51
1:B:228:PRO:HB2	1:B:249:ILE:HD12	1.93	0.51
1:A:35:ARG:HD3	1:A:198:ASP:OD2	2.10	0.51
1:B:234:ARG:O	1:B:236:PRO:HD3	2.11	0.51
1:A:116:LEU:HD22	1:A:188:MET:CE	2.38	0.50
1:A:18:LYS:HE3	1:B:154:VAL:O	2.11	0.50
1:A:116:LEU:HD23	1:A:239:ILE:HG21	1.94	0.49
1:A:23:GLY:O	1:A:24:THR:C	2.56	0.49
4:A:1015:HOH:O	1:B:104:PRO:HG2	2.12	0.49
1:A:83:TRP:CE2	1:A:143:LEU:HD11	2.49	0.47
1:B:92:PRO:O	1:B:141:MET:HE2	2.14	0.47
1:B:233:LYS:HD2	1:B:248:GLU:HG2	1.97	0.47
1:B:115:VAL:HG12	1:B:188:MET:HE1	1.96	0.47
1:B:190:GLN:HE22	1:B:232:ILE:HG21	1.79	0.47
1:A:52:LEU:HD22	1:A:249:ILE:CG1	2.44	0.47
1:A:5:LEU:HD22	1:A:220:LEU:HD23	1.97	0.47
1:B:116:LEU:HD22	4:B:859:HOH:O	2.16	0.45
1:B:225:ARG:HD3	1:B:253:ASP:O	2.17	0.45
1:A:77:VAL:HA	4:A:885:HOH:O	2.16	0.45
1:A:255:HIS:HB3	1:A:256:PRO:CD	2.46	0.45
1:B:116:LEU:CD1	1:B:188:MET:HE3	2.47	0.45
1:A:125:SER:CB	1:B:21:ARG:HH21	2.30	0.45
1:A:20:ASP:CB	1:B:126:ARG:HH12	2.20	0.45
1:B:93:VAL:O	1:B:94:TYR:C	2.59	0.45
1:B:233:LYS:HD2	1:B:248:GLU:CG	2.47	0.45
1:A:93:VAL:O	1:A:94:TYR:C	2.60	0.44
1:A:153:TYR:CE2	1:B:18:LYS:HE3	2.53	0.44
1:A:8:MET:HE1	1:A:219:GLN:NE2	2.33	0.43
1:B:51:HIS:HB3	4:B:994:HOH:O	2.18	0.42
1:A:18:LYS:HE2	1:B:124:ASP:OD1	2.19	0.42
1:A:153:TYR:CZ	1:B:18:LYS:HE3	2.55	0.41
1:A:184:LEU:O	1:A:188:MET:HG3	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:260:ALA:HB1	1:A:261:PRO:HD2	2.01	0.41
1:A:102:PRO:HG2	4:A:1008:HOH:O	2.20	0.41
1:B:111:GLN:O	1:B:115:VAL:HG23	2.20	0.41
1:A:128:ILE:CG2	1:A:152:PHE:HB2	2.51	0.41
1:B:116:LEU:HD13	1:B:239:ILE:HG21	2.03	0.41
1:A:115:VAL:HG12	1:A:188:MET:HE1	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	261/264 (99%)	252 (97%)	8 (3%)	1 (0%)	30	27
1	B	261/264 (99%)	255 (98%)	5 (2%)	1 (0%)	30	27
All	All	522/528 (99%)	507 (97%)	13 (2%)	2 (0%)	30	27

All (2) Ramachandran outliers are listed below:

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

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	232/232 (100%)	224 (97%)	8 (3%)	32	33
1	B	232/232 (100%)	227 (98%)	5 (2%)	45	50
All	All	464/464 (100%)	451 (97%)	13 (3%)	38	41

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	14	GLU
1	A	18	LYS
1	A	20	ASP
1	A	22	THR
1	A	27	LEU
1	A	115	VAL
1	A	116	LEU
1	A	249	ILE
1	B	73	HIS
1	B	116	LEU
1	B	172	LEU
1	B	231	ILE
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	64	GLN
1	A	117	ASN
1	A	219	GLN
1	B	87	ASN
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	B	1	1	9,10,11	1.84	3 (33%)	9,11,13	3.55	5 (55%)
1	CXM	A	1	1	9,10,11	2.09	2 (22%)	9,11,13	3.62	4 (44%)

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

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	1	CXM	CA-N	-5.08	1.39	1.46
1	B	1	CXM	CA-N	-3.86	1.41	1.46
1	B	1	CXM	ON1-CN	2.94	1.27	1.21
1	A	1	CXM	ON1-CN	2.29	1.25	1.21
1	B	1	CXM	O-C	2.01	1.27	1.20

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1	CXM	CA-N-CN	-8.60	106.33	122.11
1	B	1	CXM	ON2-CN-ON1	-6.20	74.29	121.85
1	B	1	CXM	CA-N-CN	-6.11	110.90	122.11
1	A	1	CXM	ON2-CN-ON1	-4.42	87.95	121.85
1	A	1	CXM	CB-CA-N	3.82	117.47	110.52
1	B	1	CXM	CB-CA-N	3.72	117.30	110.52
1	B	1	CXM	C-CA-N	-3.47	102.80	109.50
1	B	1	CXM	O-C-CA	-2.96	117.15	124.77
1	A	1	CXM	C-CA-N	-2.08	105.49	109.50

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	1	CXM	ON2-CN-N-CA
1	B	1	CXM	ON2-CN-N-CA
1	A	1	CXM	ON1-CN-N-CA
1	B	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
3	TP3	B	702[B]	-	30,30,30	3.77	5 (16%)	43,43,43	2.13	15 (34%)
2	SO4	B	802	-	4,4,4	0.65	0	6,6,6	0.41	0
3	TP3	B	702[A]	-	30,30,30	3.77	5 (16%)	43,43,43	2.16	15 (34%)
2	SO4	A	811	-	4,4,4	0.68	0	6,6,6	0.16	0
2	SO4	A	801	-	4,4,4	0.63	0	6,6,6	0.31	0
3	TP3	A	701	-	30,30,30	3.83	3 (10%)	43,43,43	2.12	16 (37%)
2	SO4	A	812	-	4,4,4	0.66	0	6,6,6	0.17	0
2	SO4	B	814	-	4,4,4	0.67	0	6,6,6	0.33	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
3	TP3	B	702[A]	-	-	8/33/43/43	0/2/2/2
3	TP3	B	702[B]	-	-	7/33/43/43	0/2/2/2
3	TP3	A	701	-	-	4/33/43/43	0/2/2/2

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	701	TP3	O12-S9	15.10	1.59	1.43
3	B	702[A]	TP3	O12-S9	14.00	1.58	1.43
3	B	702[B]	TP3	O12-S9	14.00	1.58	1.43
3	B	702[A]	TP3	O11-S9	12.49	1.56	1.43
3	B	702[B]	TP3	O11-S9	12.49	1.56	1.43
3	A	701	TP3	O11-S9	12.16	1.56	1.43
3	B	702[A]	TP3	S9-N3	6.64	1.73	1.63
3	B	702[B]	TP3	S9-N3	6.64	1.73	1.63
3	A	701	TP3	S9-N3	6.11	1.72	1.63
3	B	702[A]	TP3	C2-N3	-2.19	1.45	1.48
3	B	702[B]	TP3	C2-N3	-2.19	1.45	1.48
3	B	702[A]	TP3	C20-C25	2.03	1.57	1.52
3	B	702[B]	TP3	C20-C25	2.03	1.57	1.52

All (46) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	701	TP3	C6-C4-N3	4.88	123.13	111.91
3	A	701	TP3	O12-S9-O11	-4.80	112.10	119.59
3	A	701	TP3	C25-C20-N19	-4.68	99.72	110.57
3	B	702[A]	TP3	C13-C10-S9	-4.55	115.22	119.73
3	B	702[B]	TP3	C13-C10-S9	-4.55	115.22	119.73
3	B	702[A]	TP3	C6-C4-N3	4.45	122.14	111.91
3	B	702[B]	TP3	C6-C4-N3	4.45	122.14	111.91
3	B	702[A]	TP3	O8-C6-O7	-4.24	114.46	124.08
3	B	702[B]	TP3	O8-C6-O7	-4.24	114.46	124.08
3	B	702[A]	TP3	O12-S9-O11	-3.82	113.62	119.59
3	B	702[B]	TP3	O12-S9-O11	-3.82	113.62	119.59
3	B	702[A]	TP3	C22-C21-C20	3.81	120.18	113.16
3	B	702[A]	TP3	C20-N19-C18	3.79	130.65	121.56
3	B	702[B]	TP3	C20-N19-C18	3.79	130.65	121.56
3	A	701	TP3	O11-S9-N3	3.70	113.45	106.97
3	A	701	TP3	C17-C10-C13	3.47	124.99	120.47
3	B	702[B]	TP3	C22-C21-C20	3.43	119.49	113.16

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	702[A]	TP3	C21-C20-N19	3.41	117.66	110.91
3	B	702[B]	TP3	C21-C20-N19	3.41	117.66	110.91
3	A	701	TP3	C21-C20-N19	3.40	117.63	110.91
3	B	702[A]	TP3	C1-C2-N3	-3.04	98.12	103.32
3	B	702[B]	TP3	C1-C2-N3	-3.04	98.12	103.32
3	B	702[A]	TP3	C2-N3-C4	2.89	115.22	111.25
3	B	702[B]	TP3	C2-N3-C4	2.89	115.22	111.25
3	B	702[A]	TP3	C25-C20-N19	-2.81	104.06	110.57
3	B	702[B]	TP3	C25-C20-N19	-2.81	104.06	110.57
3	B	702[A]	TP3	O7-C6-C4	2.79	131.00	122.44
3	B	702[B]	TP3	O7-C6-C4	2.79	131.00	122.44
3	B	702[A]	TP3	C17-C10-S9	2.56	122.27	119.73
3	B	702[B]	TP3	C17-C10-S9	2.56	122.27	119.73
3	A	701	TP3	C16-C17-C10	-2.51	117.00	119.44
3	A	701	TP3	C1-C2-N3	-2.49	99.06	103.32
3	A	701	TP3	O24-C18-N19	-2.47	117.76	122.47
3	A	701	TP3	C17-C10-S9	-2.40	117.35	119.73
3	B	702[A]	TP3	O11-S9-N3	2.30	110.99	106.97
3	B	702[B]	TP3	O11-S9-N3	2.30	110.99	106.97
3	A	701	TP3	O8-C6-O7	-2.29	118.89	124.08
3	B	702[A]	TP3	C5-C4-C6	2.17	115.88	111.26
3	B	702[B]	TP3	C5-C4-C6	2.17	115.88	111.26
3	A	701	TP3	C4-N3-S9	2.15	124.65	119.77
3	A	701	TP3	C2-N3-C4	2.13	114.18	111.25
3	A	701	TP3	C16-C15-C14	2.13	121.28	118.57
3	A	701	TP3	C13-C10-S9	-2.09	117.66	119.73
3	B	702[A]	TP3	C10-S9-N3	-2.09	103.71	107.36
3	B	702[B]	TP3	C10-S9-N3	-2.09	103.71	107.36
3	A	701	TP3	O26-C25-C20	-2.07	115.56	122.26

There are no chirality outliers.

All (19) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	701	TP3	N19-C20-C21-C22
3	B	702[A]	TP3	N19-C20-C21-C22
3	A	701	TP3	C25-C20-C21-C22
3	B	702[A]	TP3	C25-C20-C21-C22
3	B	702[B]	TP3	C20-C21-C22-C23
3	B	702[A]	TP3	C13-C10-S9-O12
3	B	702[B]	TP3	C13-C10-S9-O12
3	B	702[A]	TP3	C17-C10-S9-O12

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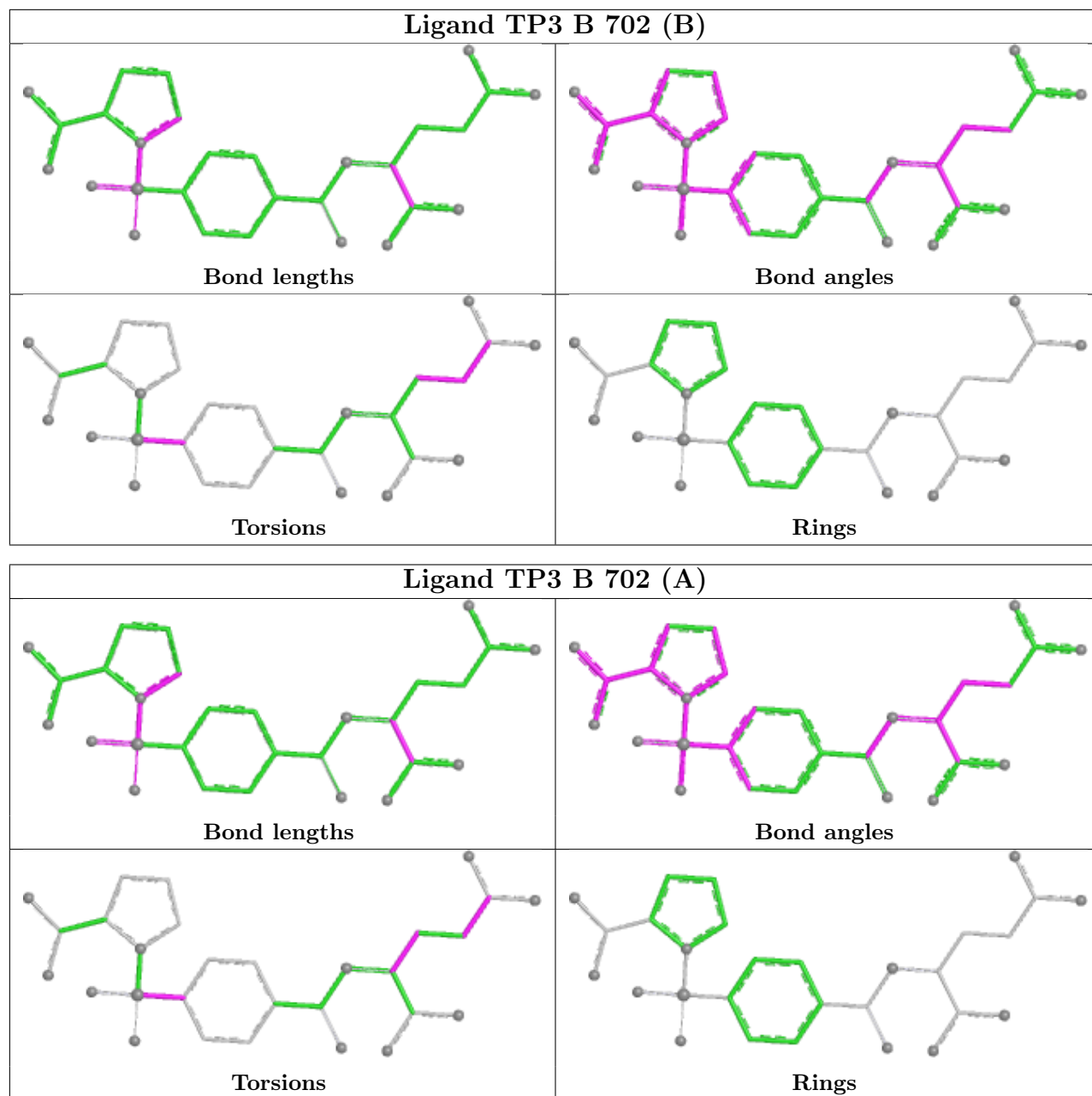
Mol	Chain	Res	Type	Atoms
3	B	702[B]	TP3	C17-C10-S9-O12
3	A	701	TP3	C21-C22-C23-O29
3	A	701	TP3	C21-C22-C23-O28
3	B	702[B]	TP3	C21-C22-C23-O29
3	B	702[A]	TP3	C21-C22-C23-O29
3	B	702[B]	TP3	C21-C22-C23-O28
3	B	702[A]	TP3	C13-C10-S9-N3
3	B	702[B]	TP3	C13-C10-S9-N3
3	B	702[A]	TP3	C21-C22-C23-O28
3	B	702[A]	TP3	C17-C10-S9-N3
3	B	702[B]	TP3	C17-C10-S9-N3

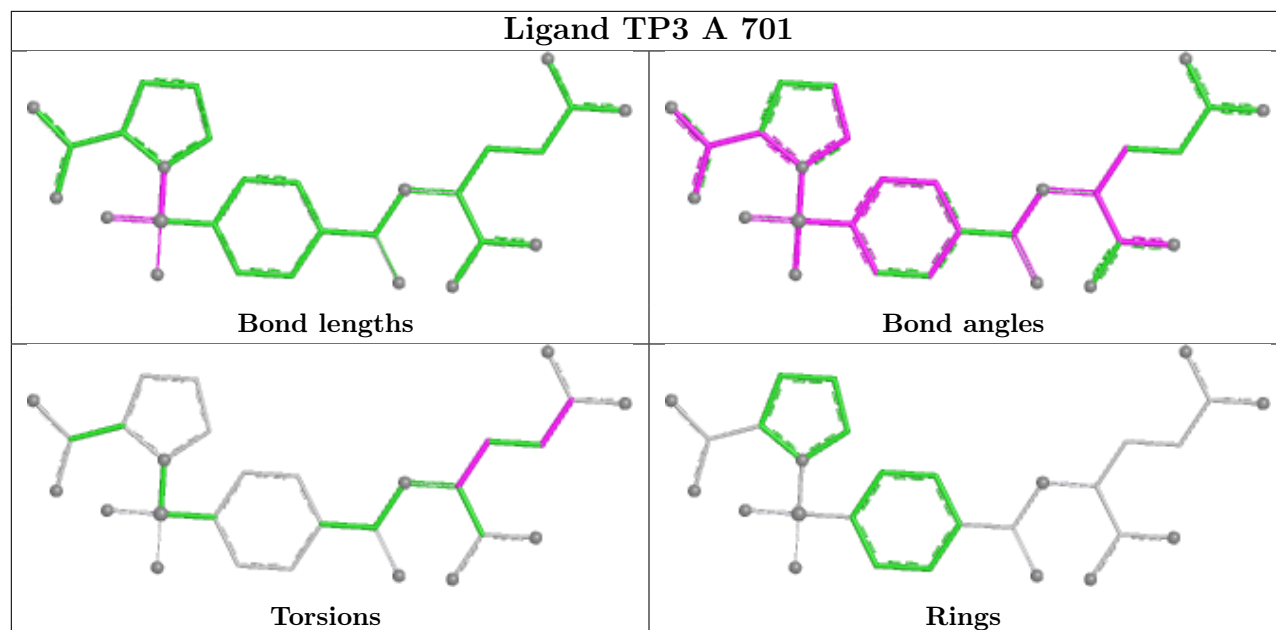
There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	702[B]	TP3	1	0
3	A	701	TP3	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.