Full wwPDB X-ray Structure Validation Report ⓘ

Feb 3, 2024 – 01:26 PM EST

PDB ID	:	1LCE
Title	:	LACTOBACILLUS CASEI THYMIDYLATE SYNTHASE TERNARY COMPLEX WITH DUMP AND CH2THF
Authors	:	Birdsall, D.L.; Finer-Moore, J.; Stroud, R.M.
Deposited on	:	1995-06-22
Resolution	:	2.50 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:

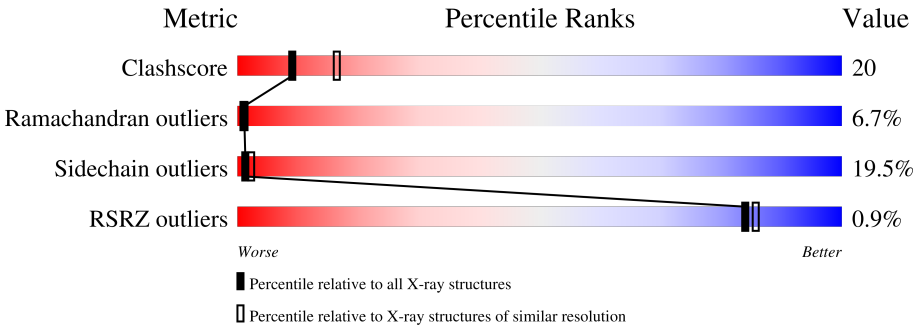
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)
RSRZ outliers	127900	4559 (2.50-2.50)

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 <=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	316	% 39% 40% 17% 5%

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
2	UMP	A	317	X	-	-	-
3	TMF	A	318	X	-	X	-

2 Entry composition

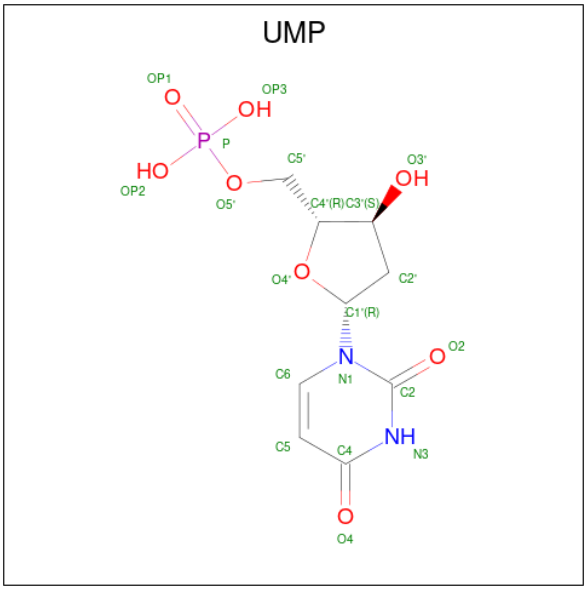
There are 4 unique types of molecules in this entry. The entry contains 3283 atoms, of which 610 are hydrogens and 0 are deuteriums.

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.

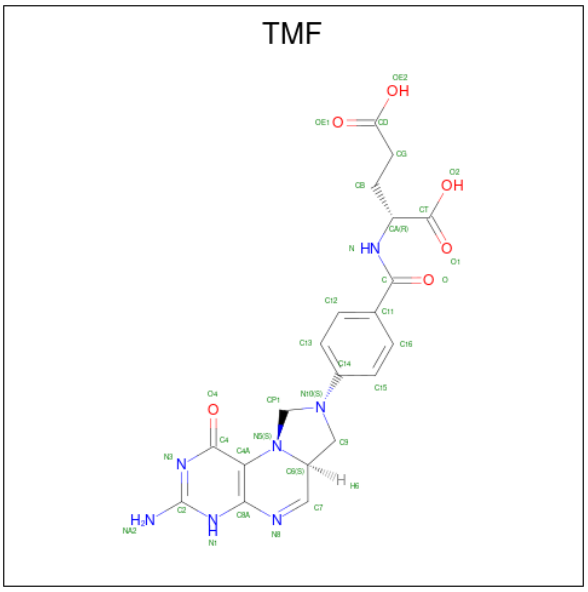
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
1	A	316	3139	1677	549	438	467	8	0	0	0

- Molecule 2 is 2'-DEOXYURIDINE 5'-MONOPHOSPHATE (three-letter code: UMP) (formula: $C_9H_{13}N_2O_8P$).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
			Total	C	H	N	O	P		
2	A	1	21	9	1	2	8	1	0	0

- Molecule 3 is 5,10-METHYLENE-6-HYDROFOLIC ACID (three-letter code: TMF) (formula: $C_{20}H_{21}N_7O_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			33	20	7	6		

- Molecule 4 is water.

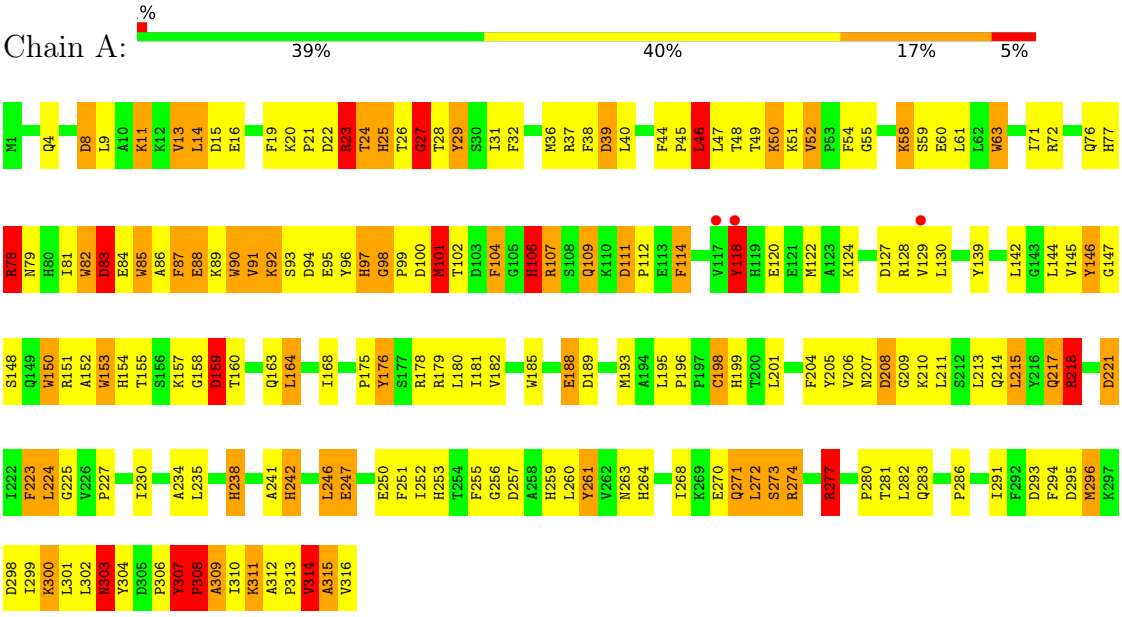
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	30	Total	H	O	0	0
			90	60	30		

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: THYMIDYLATE SYNTHASE

Chain A:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 61 2 2	Depositor
Cell constants a, b, c, α, β, γ	78.90Å 78.90Å 229.70Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	15.00 – 2.50 58.72 – 2.43	Depositor EDS
% Data completeness (in resolution range)	58.5 (15.00-2.50) 55.5 (58.72-2.43)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.29 (at 2.42Å)	Xtriage
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.204 , (Not available) 0.204 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å^2)	22.2	Xtriage
Anisotropy	0.643	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.30 , 144.2	EDS
L-test for twinning[2]	$< \|L\| > = 0.44$, $< L^2 > = 0.27$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.89	EDS
Total number of atoms	3283	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 4.83% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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: UMP, TMF

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	1.24	4/2674 (0.1%)	2.18	117/3634 (3.2%)

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	3

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	120	GLU	CB-CG	5.57	1.62	1.52
1	A	107	ARG	NE-CZ	5.36	1.40	1.33
1	A	120	GLU	CG-CD	5.27	1.59	1.51
1	A	316	VAL	CA-CB	5.07	1.65	1.54

All (117) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	122	MET	CG-SD-CE	-10.96	82.67	100.20
1	A	296	MET	CG-SD-CE	-10.37	83.62	100.20
1	A	78	ARG	NE-CZ-NH1	10.08	125.34	120.30
1	A	122	MET	CA-CB-CG	-8.97	98.05	113.30
1	A	308	PRO	N-CA-CB	-8.88	92.64	103.30
1	A	101	MET	CA-CB-CG	-8.84	98.27	113.30
1	A	82	TRP	CD1-CG-CD2	8.77	113.31	106.30
1	A	150	TRP	CD1-CG-CD2	8.70	113.26	106.30
1	A	303	ASN	CA-C-N	-8.65	98.17	117.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	185	TRP	CD1-CG-CD2	8.29	112.93	106.30
1	A	179	ARG	NE-CZ-NH2	-7.74	116.43	120.30
1	A	153	TRP	CD1-CG-CD2	7.66	112.43	106.30
1	A	185	TRP	CE2-CD2-CG	-7.65	101.18	107.30
1	A	150	TRP	CE2-CD2-CG	-7.61	101.21	107.30
1	A	37	ARG	NE-CZ-NH2	-7.60	116.50	120.30
1	A	208	ASP	CA-C-N	-7.58	101.05	116.20
1	A	85	TRP	CD1-CG-CD2	7.53	112.32	106.30
1	A	82	TRP	CE2-CD2-CG	-7.50	101.30	107.30
1	A	77	HIS	CA-CB-CG	-7.50	100.86	113.60
1	A	97	HIS	CA-C-N	-7.48	101.23	116.20
1	A	78	ARG	CD-NE-CZ	7.45	134.02	123.60
1	A	23	ARG	CA-CB-CG	7.43	129.75	113.40
1	A	88	GLU	CA-CB-CG	7.34	129.54	113.40
1	A	15	ASP	CB-CG-OD1	7.31	124.88	118.30
1	A	124	LYS	CA-CB-CG	7.29	129.43	113.40
1	A	153	TRP	CE2-CD2-CG	-7.28	101.48	107.30
1	A	36	MET	CA-CB-CG	7.27	125.65	113.30
1	A	46	LEU	CA-CB-CG	7.12	131.67	115.30
1	A	106	HIS	CA-C-N	-7.09	101.60	117.20
1	A	224	LEU	CA-C-N	6.99	130.17	116.20
1	A	128	ARG	NE-CZ-NH2	-6.97	116.81	120.30
1	A	63	TRP	CD1-CG-CD2	6.74	111.69	106.30
1	A	27	GLY	CA-C-N	-6.67	102.51	117.20
1	A	97	HIS	N-CA-C	6.61	128.84	111.00
1	A	82	TRP	CG-CD1-NE1	-6.56	103.54	110.10
1	A	71	ILE	CA-CB-CG2	-6.54	97.82	110.90
1	A	151	ARG	NE-CZ-NH1	6.52	123.56	120.30
1	A	63	TRP	CE2-CD2-CG	-6.40	102.18	107.30
1	A	147	GLY	O-C-N	-6.33	112.57	122.70
1	A	242	HIS	CA-CB-CG	6.32	124.34	113.60
1	A	72	ARG	NE-CZ-NH1	6.32	123.46	120.30
1	A	85	TRP	CE2-CD2-CG	-6.30	102.26	107.30
1	A	90	TRP	CE2-CD2-CG	-6.22	102.32	107.30
1	A	295	ASP	CB-CG-OD1	6.20	123.88	118.30
1	A	176	TYR	CB-CG-CD1	-6.19	117.28	121.00
1	A	15	ASP	CA-CB-CG	6.10	126.82	113.40
1	A	109	GLN	CA-CB-CG	6.10	126.83	113.40
1	A	303	ASN	CA-C-O	6.09	132.89	120.10
1	A	27	GLY	CA-C-O	6.07	131.53	120.60
1	A	93	SER	CA-C-N	-6.05	103.88	117.20
1	A	157	LYS	N-CA-CB	-6.05	99.71	110.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	261	TYR	CB-CG-CD1	-5.97	117.42	121.00
1	A	193	MET	O-C-N	5.93	132.19	122.70
1	A	193	MET	CA-C-N	-5.92	104.17	117.20
1	A	114	PHE	N-CA-C	5.90	126.93	111.00
1	A	23	ARG	NE-CZ-NH2	5.89	123.25	120.30
1	A	72	ARG	NE-CZ-NH2	-5.85	117.37	120.30
1	A	277	ARG	CA-CB-CG	5.85	126.26	113.40
1	A	107	ARG	N-CA-CB	-5.84	100.08	110.60
1	A	164	LEU	CA-CB-CG	5.83	128.71	115.30
1	A	272	LEU	CA-CB-CG	-5.81	101.94	115.30
1	A	298	ASP	CB-CG-OD1	5.79	123.52	118.30
1	A	88	GLU	N-CA-CB	5.77	120.98	110.60
1	A	106	HIS	N-CA-C	5.74	126.48	111.00
1	A	198	CYS	CA-CB-SG	5.72	124.30	114.00
1	A	58	LYS	CA-CB-CG	5.72	125.98	113.40
1	A	124	LYS	CD-CE-NZ	5.71	124.84	111.70
1	A	8	ASP	CA-CB-CG	5.70	125.94	113.40
1	A	9	LEU	CA-CB-CG	5.65	128.30	115.30
1	A	146	TYR	CA-C-N	5.60	127.40	116.20
1	A	201	LEU	CA-CB-CG	5.58	128.15	115.30
1	A	89	LYS	CA-C-N	5.58	129.47	117.20
1	A	92	LYS	CA-CB-CG	5.55	125.61	113.40
1	A	13	VAL	CG1-CB-CG2	5.54	119.76	110.90
1	A	217	GLN	N-CA-C	-5.54	96.05	111.00
1	A	198	CYS	O-C-N	-5.54	113.84	122.70
1	A	218	ARG	NE-CZ-NH1	5.53	123.07	120.30
1	A	129	VAL	CG1-CB-CG2	-5.52	102.07	110.90
1	A	247	GLU	CA-CB-CG	5.50	125.49	113.40
1	A	150	TRP	CG-CD1-NE1	-5.49	104.61	110.10
1	A	309	ALA	CA-C-N	5.49	129.28	117.20
1	A	291	ILE	CA-CB-CG2	-5.49	99.92	110.90
1	A	15	ASP	CB-CG-OD2	-5.47	113.38	118.30
1	A	250	GLU	N-CA-CB	-5.46	100.77	110.60
1	A	178	ARG	N-CA-CB	-5.46	100.78	110.60
1	A	87	PHE	CB-CG-CD2	-5.45	116.99	120.80
1	A	277	ARG	NE-CZ-NH1	5.45	123.03	120.30
1	A	84	GLU	CA-CB-CG	5.38	125.24	113.40
1	A	85	TRP	CG-CD1-NE1	-5.37	104.73	110.10
1	A	83	ASP	N-CA-C	5.34	125.43	111.00
1	A	128	ARG	NE-CZ-NH1	5.34	122.97	120.30
1	A	40	LEU	CA-C-N	5.32	128.89	117.20
1	A	308	PRO	N-CD-CG	-5.30	95.25	103.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	29	TYR	N-CA-C	-5.28	96.74	111.00
1	A	71	ILE	CA-CB-CG1	5.26	120.99	111.00
1	A	14	LEU	N-CA-C	-5.24	96.85	111.00
1	A	159	ASP	C-N-CA	5.24	134.80	121.70
1	A	88	GLU	CB-CA-C	-5.24	99.93	110.40
1	A	188	GLU	N-CA-CB	5.23	120.02	110.60
1	A	209	GLY	N-CA-C	-5.23	100.02	113.10
1	A	52	VAL	CG1-CB-CG2	-5.22	102.54	110.90
1	A	160	THR	N-CA-CB	-5.22	100.38	110.30
1	A	118	TYR	CG-CD2-CE2	-5.21	117.13	121.30
1	A	120	GLU	CA-CB-CG	5.17	124.78	113.40
1	A	188	GLU	CA-CB-CG	5.17	124.77	113.40
1	A	29	TYR	CG-CD2-CE2	-5.16	117.17	121.30
1	A	13	VAL	CA-CB-CG1	-5.16	103.16	110.90
1	A	97	HIS	CA-CB-CG	-5.12	104.90	113.60
1	A	146	TYR	CB-CG-CD2	-5.11	117.94	121.00
1	A	221	ASP	CB-CG-OD1	5.11	122.90	118.30
1	A	39	ASP	N-CA-C	-5.10	97.23	111.00
1	A	251	PHE	N-CA-C	-5.06	97.35	111.00
1	A	144	LEU	CA-C-N	-5.05	106.08	117.20
1	A	188	GLU	CA-C-N	-5.03	106.14	117.20
1	A	79	ASN	O-C-N	-5.02	114.67	122.70
1	A	189	ASP	CB-CG-OD1	5.01	122.81	118.30
1	A	25	HIS	N-CA-C	5.00	124.51	111.00

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	111	ASP	Peptide
1	A	307	TYR	Peptide
1	A	98	GLY	Peptide

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	2590	549	2496	100	0
2	A	20	1	10	4	0
3	A	33	0	19	17	0
4	A	30	60	0	1	0
All	All	2673	610	2525	104	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 20.

All (104) 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:8:ASP:HA	1:A:11:LYS:HE3	1.54	0.90
1:A:81:ILE:HG21	3:A:318:TMF:H92	1.57	0.86
1:A:199:HIS:HB3	1:A:215:LEU:HD21	1.60	0.82
1:A:211:LEU:HB2	1:A:246:LEU:HD13	1.61	0.80
1:A:85:TRP:HA	1:A:88:GLU:HG2	1.65	0.77
1:A:101:MET:SD	1:A:104:PHE:HB3	2.30	0.72
1:A:195:LEU:HG	3:A:318:TMF:NA2	2.06	0.71
1:A:54:PHE:CD2	1:A:301:LEU:HD22	2.29	0.68
1:A:51:LYS:HD2	1:A:306:PRO:HG3	1.76	0.66
1:A:104:PHE:HB2	1:A:118:TYR:CD1	2.31	0.66
1:A:175:PRO:HB3	1:A:206:VAL:HG11	1.77	0.65
1:A:82:TRP:CZ2	3:A:318:TMF:H6	2.32	0.65
1:A:195:LEU:HG	3:A:318:TMF:HN21	1.61	0.65
1:A:20:LYS:HB3	1:A:28:THR:OG1	1.96	0.65
1:A:95:GLU:HB2	1:A:139:TYR:OH	1.96	0.64
1:A:55:GLY:HA2	1:A:58:LYS:HD2	1.78	0.64
1:A:274:ARG:HB2	1:A:307:TYR:CE2	2.33	0.64
2:A:317:UMP:O4	3:A:318:TMF:HCP1	1.97	0.64
1:A:204:PHE:CD2	1:A:211:LEU:HD21	2.34	0.63
1:A:90:TRP:HZ3	1:A:95:GLU:HB3	1.62	0.62
1:A:195:LEU:CG	3:A:318:TMF:HN21	2.14	0.60
1:A:235:LEU:HD11	1:A:299:ILE:HG21	1.84	0.59
1:A:270:GLU:O	1:A:273:SER:HB3	2.02	0.58
1:A:48:THR:HB	1:A:277:ARG:HG3	1.86	0.58
1:A:48:THR:O	1:A:306:PRO:HA	2.04	0.58
1:A:49:THR:O	1:A:50:LYS:HB3	2.05	0.57
1:A:205:TYR:HE2	1:A:207:ASN:OD1	1.87	0.57
1:A:81:ILE:CG2	3:A:318:TMF:H92	2.34	0.55
1:A:224:LEU:HA	3:A:318:TMF:CT	2.37	0.55
1:A:309:ALA:O	1:A:310:ILE:HG13	2.07	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:19:PHE:HA	1:A:28:THR:O	2.09	0.53
1:A:85:TRP:O	1:A:88:GLU:HB2	2.09	0.53
1:A:46:LEU:HD12	1:A:280:PRO:HG3	1.91	0.53
1:A:164:LEU:O	1:A:168:ILE:HG12	2.10	0.52
1:A:224:LEU:HD21	1:A:312:ALA:HB3	1.91	0.52
3:A:318:TMF:H15	3:A:318:TMF:N8	2.24	0.52
1:A:283:GLN:HE21	1:A:300:LYS:NZ	2.08	0.51
1:A:281:THR:HB	1:A:302:LEU:HB2	1.92	0.51
1:A:224:LEU:HA	3:A:318:TMF:O2	2.11	0.50
1:A:230:ILE:HG12	1:A:253:HIS:ND1	2.25	0.50
1:A:83:ASP:O	1:A:86:ALA:HB3	2.11	0.50
3:A:318:TMF:H15	3:A:318:TMF:C7	2.42	0.50
1:A:38:PHE:CE2	1:A:230:ILE:HD13	2.46	0.49
1:A:198:CYS:O	1:A:217:GLN:HA	2.11	0.49
1:A:20:LYS:O	1:A:27:GLY:HA2	2.13	0.49
1:A:90:TRP:CZ3	1:A:95:GLU:HB3	2.45	0.48
1:A:204:PHE:HB3	1:A:211:LEU:HD11	1.94	0.48
1:A:195:LEU:HD12	1:A:196:PRO:HD2	1.94	0.48
1:A:23:ARG:NH1	2:A:317:UMP:H5'	2.28	0.48
1:A:307:TYR:HB3	1:A:308:PRO:HB2	1.95	0.48
1:A:224:LEU:HD22	3:A:318:TMF:HG2	1.95	0.47
1:A:283:GLN:NE2	1:A:300:LYS:HZ3	2.12	0.47
1:A:205:TYR:CE2	1:A:207:ASN:HA	2.49	0.47
1:A:29:TYR:O	1:A:259:HIS:HA	2.14	0.47
1:A:44:PHE:HB3	1:A:280:PRO:HG2	1.97	0.47
1:A:47:LEU:HD22	1:A:227:PRO:HB3	1.97	0.47
1:A:32:PHE:HA	1:A:256:GLY:O	2.15	0.47
1:A:264:HIS:CD2	1:A:312:ALA:HB1	2.50	0.47
1:A:47:LEU:O	1:A:304:TYR:HE1	1.98	0.46
1:A:223:PHE:CE2	1:A:312:ALA:HB2	2.51	0.46
1:A:274:ARG:HB2	1:A:307:TYR:CZ	2.50	0.46
1:A:60:GLU:O	1:A:63:TRP:HB3	2.16	0.46
1:A:54:PHE:CG	1:A:301:LEU:HD22	2.51	0.45
1:A:198:CYS:SG	1:A:218:ARG:NH2	2.83	0.45
1:A:204:PHE:CE2	1:A:213:LEU:HD12	2.51	0.45
1:A:214:GLN:HG3	1:A:252:ILE:HG22	1.99	0.45
1:A:98:GLY:HA3	1:A:99:PRO:HD3	1.52	0.45
1:A:221:ASP:HB2	2:A:317:UMP:H2"	1.97	0.45
1:A:104:PHE:HA	1:A:114:PHE:CD2	2.52	0.45
1:A:257:ASP:OD1	1:A:259:HIS:ND1	2.50	0.45
1:A:106:HIS:NE2	1:A:314:VAL:HB	2.32	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:195:LEU:CG	3:A:318:TMF:NA2	2.76	0.45
1:A:24:THR:O	1:A:26:THR:N	2.50	0.45
1:A:85:TRP:CH2	1:A:195:LEU:HB3	2.51	0.45
1:A:91:VAL:HA	1:A:96:TYR:CD1	2.52	0.45
1:A:106:HIS:HE1	1:A:315:ALA:O	2.00	0.45
1:A:224:LEU:HD22	3:A:318:TMF:CG	2.47	0.45
1:A:242:HIS:HE1	1:A:286:PRO:HA	1.82	0.45
1:A:44:PHE:CD1	1:A:45:PRO:HD2	2.52	0.44
1:A:234:ALA:O	1:A:238:HIS:HB2	2.18	0.44
1:A:274:ARG:CB	1:A:307:TYR:CE2	3.00	0.44
1:A:46:LEU:HA	4:A:355:HOH:O	2.17	0.44
1:A:311:LYS:NZ	1:A:313:PRO:HD3	2.33	0.44
1:A:181:ILE:HD13	1:A:181:ILE:HG21	1.84	0.43
1:A:101:MET:HA	1:A:101:MET:CE	2.42	0.43
1:A:158:GLY:O	1:A:159:ASP:HB3	2.18	0.43
1:A:14:LEU:HD22	1:A:268:ILE:HG23	2.01	0.43
1:A:81:ILE:HG21	1:A:81:ILE:HD13	1.82	0.43
1:A:204:PHE:HD2	1:A:211:LEU:HD21	1.78	0.43
1:A:283:GLN:HE21	1:A:300:LYS:HZ3	1.66	0.43
1:A:271:GLN:HA	1:A:274:ARG:HD3	2.00	0.43
1:A:225:GLY:N	3:A:318:TMF:O1	2.52	0.42
1:A:163:GLN:HB3	1:A:182:VAL:HG13	2.00	0.42
1:A:306:PRO:HG2	1:A:307:TYR:O	2.19	0.42
1:A:90:TRP:CH2	1:A:96:TYR:HA	2.55	0.42
1:A:260:LEU:HD21	1:A:268:ILE:HG21	2.01	0.42
1:A:241:ALA:HA	1:A:246:LEU:HD12	2.02	0.41
1:A:146:TYR:O	1:A:150:TRP:HB2	2.20	0.41
1:A:224:LEU:HD22	3:A:318:TMF:CB	2.51	0.41
1:A:87:PHE:HD2	1:A:104:PHE:CE2	2.39	0.41
1:A:261:TYR:HB2	1:A:264:HIS:ND1	2.36	0.41
1:A:59:SER:HB3	1:A:296:MET:HE1	2.03	0.41
1:A:198:CYS:O	1:A:217:GLN:HG3	2.21	0.40
2:A:317:UMP:C4	3:A:318:TMF:HCP1	2.56	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	314/316 (99%)	255 (81%)	38 (12%)	21 (7%)	1 1

All (21) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	16	GLU
1	A	25	HIS
1	A	50	LYS
1	A	78	ARG
1	A	97	HIS
1	A	104	PHE
1	A	111	ASP
1	A	112	PRO
1	A	208	ASP
1	A	223	PHE
1	A	303	ASN
1	A	308	PRO
1	A	314	VAL
1	A	27	GLY
1	A	100	ASP
1	A	273	SER
1	A	315	ALA
1	A	152	ALA
1	A	210	LYS
1	A	180	LEU
1	A	307	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	277/278 (100%)	223 (80%)	54 (20%)	1 2

All (54) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	GLN
1	A	11	LYS
1	A	13	VAL
1	A	21	PRO
1	A	22	ASP
1	A	23	ARG
1	A	24	THR
1	A	31	ILE
1	A	39	ASP
1	A	46	LEU
1	A	52	VAL
1	A	61	LEU
1	A	76	GLN
1	A	78	ARG
1	A	83	ASP
1	A	91	VAL
1	A	92	LYS
1	A	94	ASP
1	A	101	MET
1	A	102	THR
1	A	106	HIS
1	A	107	ARG
1	A	109	GLN
1	A	118	TYR
1	A	127	ASP
1	A	130	LEU
1	A	142	LEU
1	A	145	VAL
1	A	148	SER
1	A	153	TRP
1	A	154	HIS
1	A	155	THR
1	A	159	ASP
1	A	176	TYR
1	A	188	GLU

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Mol	Chain	Res	Type
1	A	215	LEU
1	A	218	ARG
1	A	238	HIS
1	A	246	LEU
1	A	247	GLU
1	A	255	PHE
1	A	263	ASN
1	A	271	GLN
1	A	272	LEU
1	A	274	ARG
1	A	277	ARG
1	A	282	LEU
1	A	293	ASP
1	A	294	PHE
1	A	300	LYS
1	A	303	ASN
1	A	308	PRO
1	A	311	LYS
1	A	314	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	35	GLN
1	A	106	HIS
1	A	214	GLN
1	A	283	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 monosaccharides in this entry.

5.6 Ligand geometry (i)

2 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	TMF	A	318	-	34,36,36	3.73	15 (44%)	36,52,52	2.89	15 (41%)
2	UMP	A	317	-	21,21,21	1.58	5 (23%)	31,31,31	3.15	9 (29%)

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	TMF	A	318	-	1/1/7/9	11/21/42/42	0/3/4/4
2	UMP	A	317	-	1/1/4/4	5/10/22/22	0/2/2/2

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	318	TMF	C7-N8	15.64	1.46	1.27
3	A	318	TMF	O4-C4	9.38	1.48	1.24
3	A	318	TMF	C6-N5	4.62	1.52	1.48
3	A	318	TMF	CA-N	3.68	1.53	1.45
3	A	318	TMF	C-N	3.58	1.42	1.34
3	A	318	TMF	C2-N1	3.54	1.41	1.35
3	A	318	TMF	C4A-N5	-3.50	1.31	1.40
2	A	317	UMP	O4'-C1'	3.34	1.49	1.42
2	A	317	UMP	C2-N1	3.30	1.43	1.38
3	A	318	TMF	C4A-C4	2.50	1.44	1.41
2	A	317	UMP	C6-N1	-2.47	1.32	1.38
3	A	318	TMF	C8A-N1	-2.45	1.31	1.34
3	A	318	TMF	C4-N3	2.45	1.37	1.33
3	A	318	TMF	CP1-N5	2.35	1.51	1.45
3	A	318	TMF	C14-N10	-2.34	1.32	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	318	TMF	O2-CT	-2.33	1.22	1.30
2	A	317	UMP	C4-N3	2.20	1.42	1.38
3	A	318	TMF	CA-CT	2.20	1.58	1.52
3	A	318	TMF	C2-N3	2.06	1.39	1.35
2	A	317	UMP	P-OP2	-2.03	1.47	1.54

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	317	UMP	O4'-C1'-N1	9.16	124.23	107.86
2	A	317	UMP	C1'-N1-C6	-7.27	107.19	121.55
2	A	317	UMP	C2'-C1'-N1	7.26	130.48	113.77
3	A	318	TMF	NA2-C2-N1	6.80	127.83	117.25
2	A	317	UMP	C1'-N1-C2	6.35	130.14	117.64
3	A	318	TMF	CP1-N5-C4A	-5.76	106.39	122.96
3	A	318	TMF	C4A-C4-N3	-5.72	111.38	123.14
3	A	318	TMF	C15-C16-C11	-4.84	115.15	120.78
2	A	317	UMP	O4-C4-N3	4.73	126.25	119.31
3	A	318	TMF	CA-N-C	4.63	132.94	121.60
3	A	318	TMF	N1-C2-N3	-4.22	118.79	125.42
3	A	318	TMF	C13-C14-N10	-4.20	115.59	121.38
2	A	317	UMP	O4-C4-C5	-4.14	117.88	125.16
3	A	318	TMF	C12-C13-C14	-3.78	115.35	120.32
3	A	318	TMF	O2-CT-O1	-3.36	116.46	124.09
3	A	318	TMF	CT-CA-N	3.35	118.48	110.55
3	A	318	TMF	N1-C8A-N8	3.24	123.78	116.00
2	A	317	UMP	O2-C2-N1	3.02	126.80	122.79
3	A	318	TMF	O2-CT-CA	2.68	122.32	113.40
3	A	318	TMF	C13-C12-C11	2.67	123.89	120.78
3	A	318	TMF	NA2-C2-N3	-2.52	113.33	117.25
2	A	317	UMP	O5'-C5'-C4'	2.45	117.42	108.99
3	A	318	TMF	OE1-CD-CG	-2.42	115.30	123.08
2	A	317	UMP	O4'-C1'-C2'	-2.26	101.99	106.25

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	A	317	UMP	C1'
3	A	318	TMF	CA

All (16) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	317	UMP	C5'-O5'-P-OP1
2	A	317	UMP	C5'-O5'-P-OP2
2	A	317	UMP	C5'-O5'-P-OP3
3	A	318	TMF	C13-C14-N10-C9
3	A	318	TMF	C15-C14-N10-C9
3	A	318	TMF	C13-C14-N10-CP1
3	A	318	TMF	C15-C14-N10-CP1
3	A	318	TMF	C11-C-N-CA
3	A	318	TMF	O-C-N-CA
3	A	318	TMF	CT-CA-CB-CG
2	A	317	UMP	C2'-C1'-N1-C2
3	A	318	TMF	CB-CA-CT-O2
3	A	318	TMF	CB-CA-CT-O1
3	A	318	TMF	OE1-CD-CG-CB
2	A	317	UMP	C2'-C1'-N1-C6
3	A	318	TMF	OE2-CD-CG-CB

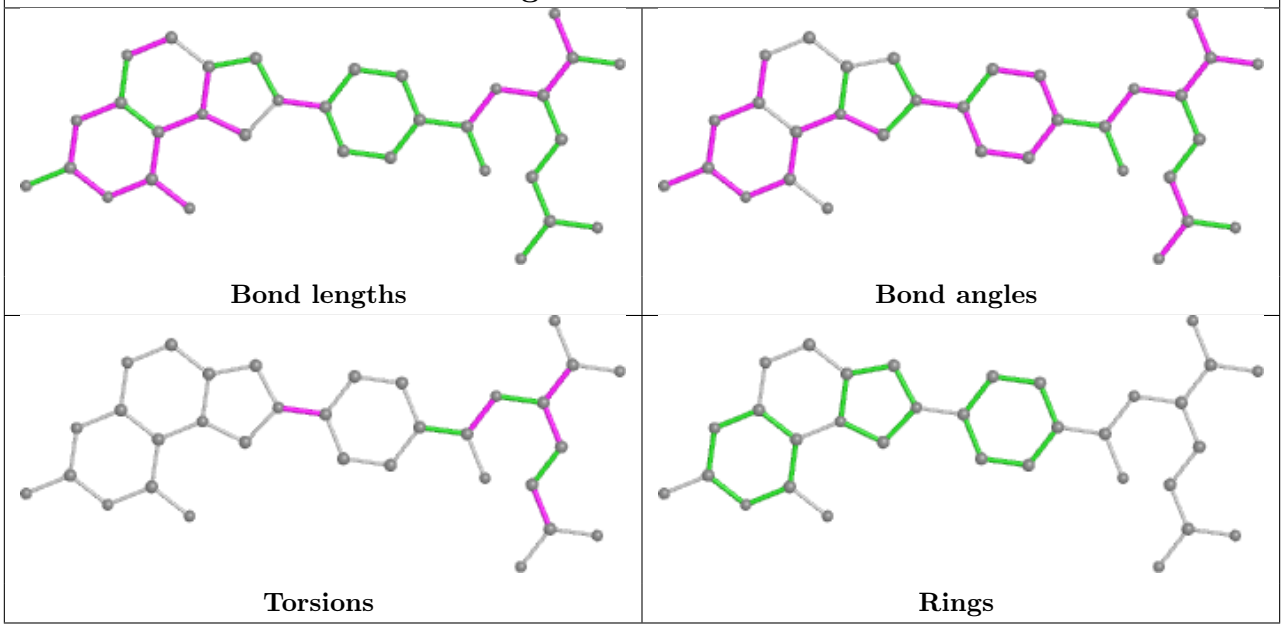
There are no ring outliers.

2 monomers are involved in 19 short contacts:

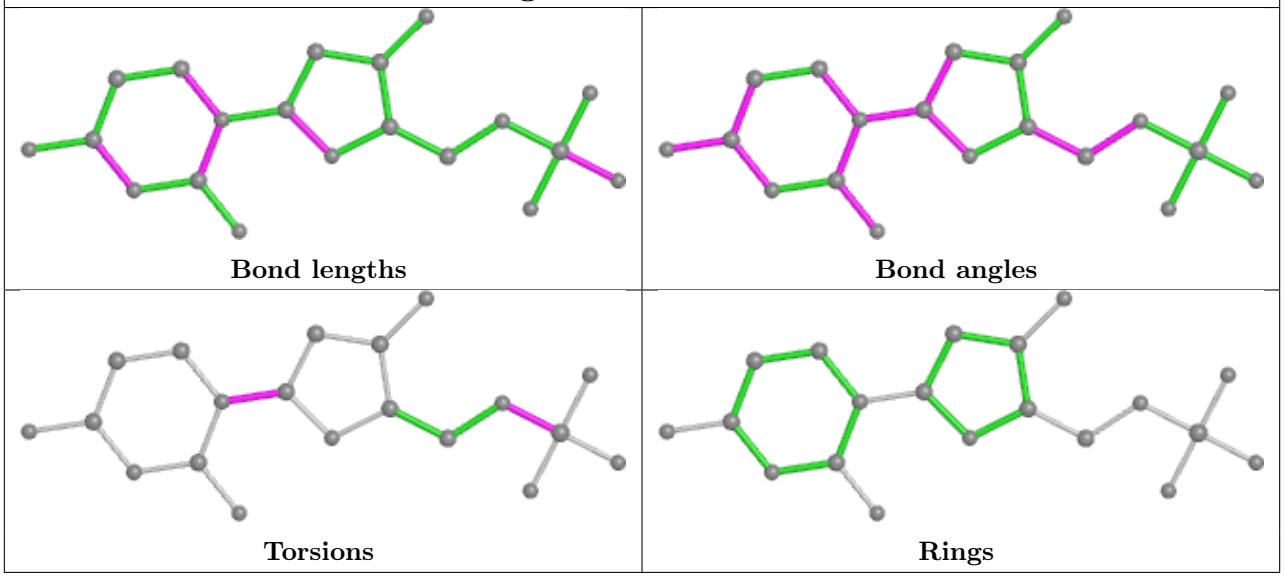
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	318	TMF	17	0
2	A	317	UMP	4	0

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 then 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.

Ligand TMF A 318



Ligand UMP A 317



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, 95^{th} 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	316/316 (100%)	-0.79	3 (0%)	84	86	2, 12, 24, 39	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	129	VAL	2.5
1	A	117	VAL	2.2
1	A	118	TYR	2.1

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

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

6.3 Carbohydrates

There are no monosaccharides in this entry.

6.4 Ligands

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, 95^{th} 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.

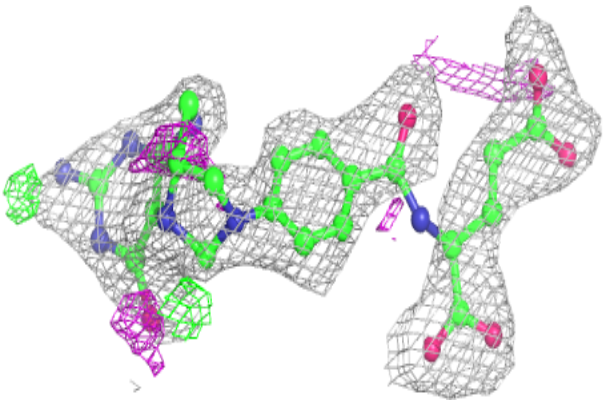
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	TMF	A	318	33/33	0.87	0.21	26,26,58,58	0
2	UMP	A	317	20/20	0.95	0.14	11,18,18,18	0

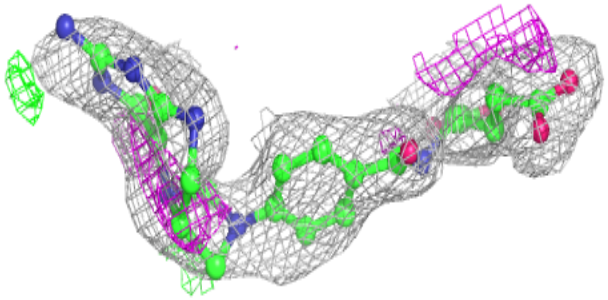
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers

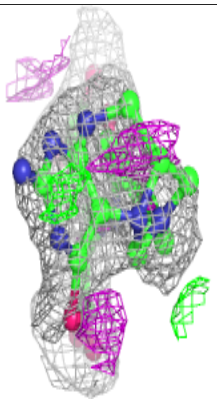
as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around TMF A 318:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

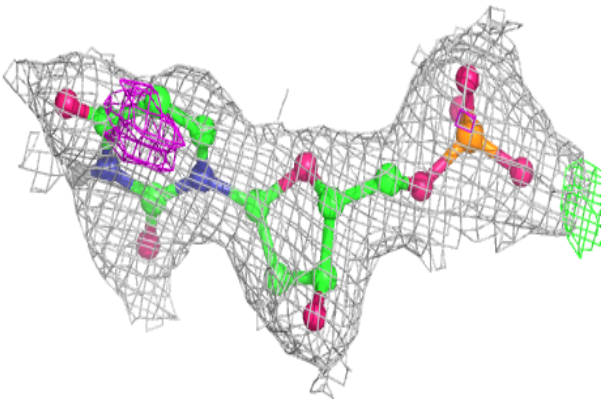


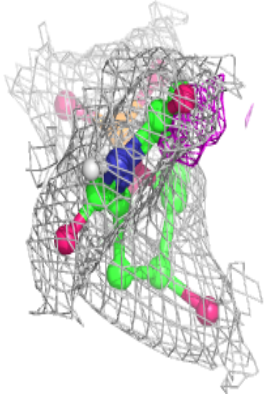




Electron density around UMP A 317:

$2mF_o$-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 (i)

There are no such residues in this entry.