



Full wwPDB X-ray Structure Validation Report ⓘ

May 16, 2020 – 04:11 pm BST

PDB ID : 1E2L
Title : Kinetics and crystal structure of the wild-type and the engineered Y101F mutant of Herpes simplex virus type 1 thymidine kinase interacting with (North)-methanocarpa-thymidine
Authors : Vogt, J.; Scapozza, L.; Schulz, G.E.
Deposited on : 2000-05-23
Resolution : 2.40 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 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 : 2.11
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) : 2.11

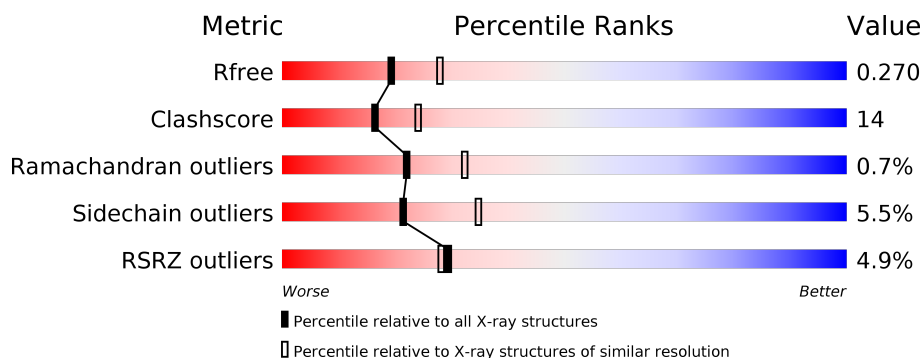
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	3907 (2.40-2.40)
Clashscore	141614	4398 (2.40-2.40)
Ramachandran outliers	138981	4318 (2.40-2.40)
Sidechain outliers	138945	4319 (2.40-2.40)
RSRZ outliers	127900	3811 (2.40-2.40)

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 ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	331	<div> <div>4%</div> <div> <div></div> <div>63%</div> <div>24%</div> <div>• 8%</div> </div> </div>
1	B	331	<div> <div>5%</div> <div> <div></div> <div>63%</div> <div>27%</div> <div>• • 6%</div> </div> </div>

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 4880 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.

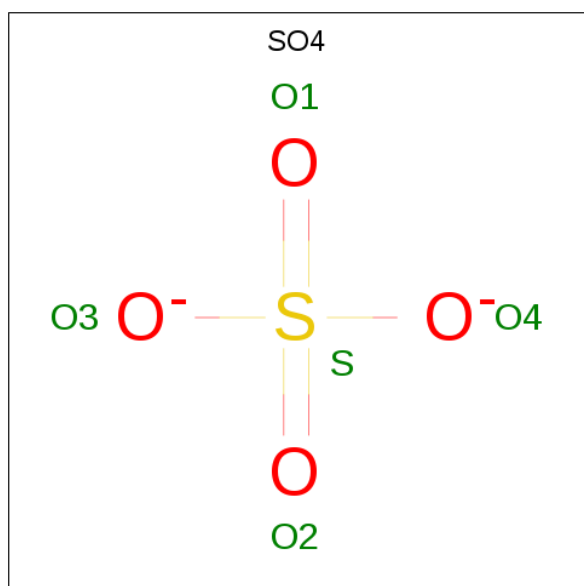
- Molecule 1 is a protein called THYMIDINE KINASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	304	Total	C	N	O	S	0	0	0
			2327	1485	406	420	16			
1	B	310	Total	C	N	O	S	0	0	0
			2357	1502	410	429	16			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	101	PHE	TYR	engineered mutation	UNP P03176
B	101	PHE	TYR	engineered mutation	UNP P03176

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



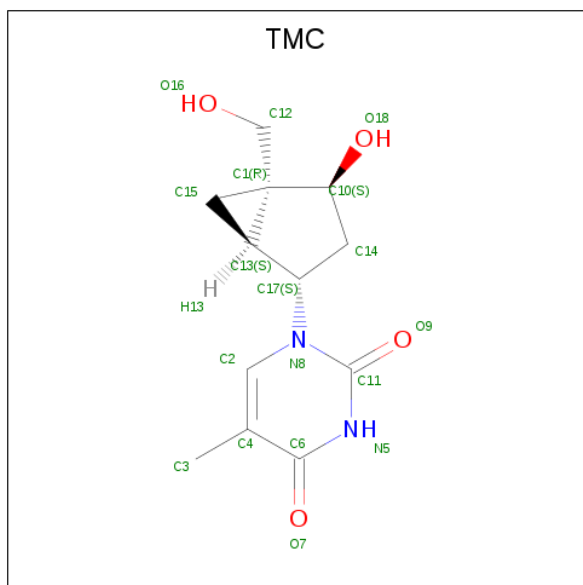
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		

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

- Molecule 3 is 1-[4-HYDROXY-5-(HYDROXYMETHYL)BICYCLO[3.1.0]HEX-2-YL]-5-METHYLPYRIMIDINE-2,4(1H,3H)-DIONE (three-letter code: TMC) (formula: C₁₂H₁₆N₂O₄).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			18	12	2	4		
3	B	1	Total	C	N	O	0	0
			18	12	2	4		

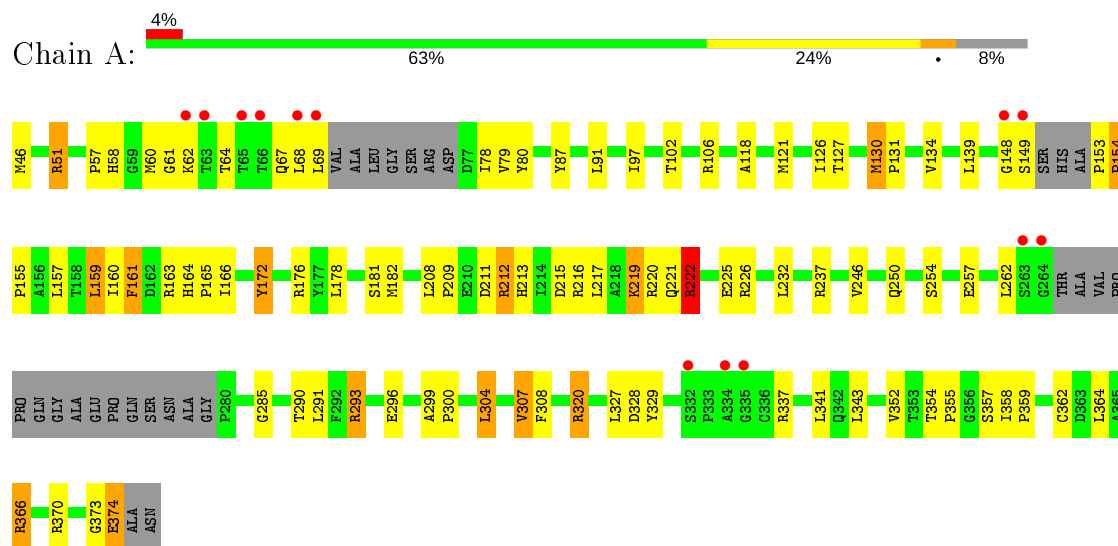
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	77	Total	O	0	0
			77	77		
4	B	73	Total	O	0	0
			73	73		

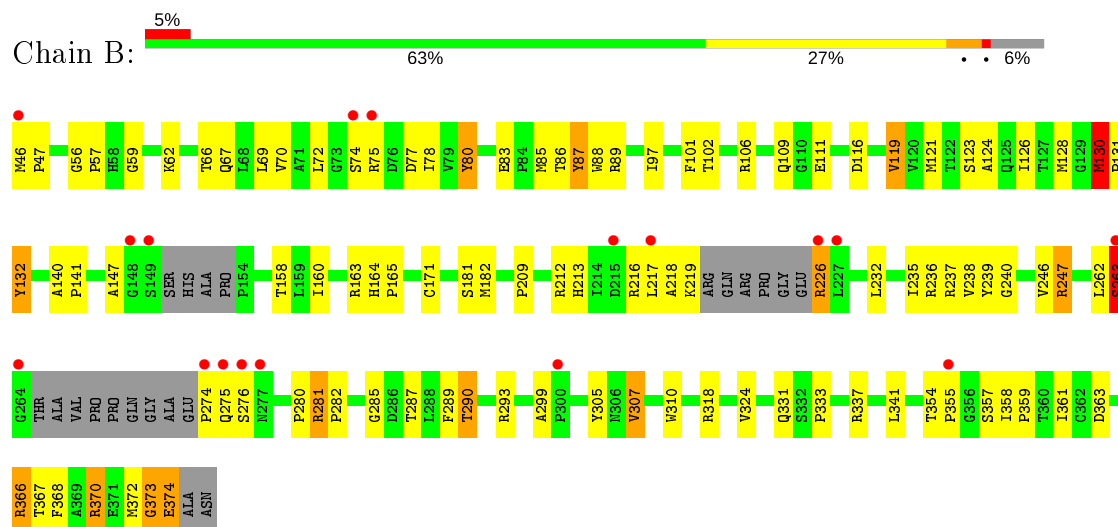
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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: THYMIDINE KINASE



• Molecule 1: THYMIDINE KINASE



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	114.00Å 118.30Å 108.60Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.40 24.99 – 2.30	Depositor EDS
% Data completeness (in resolution range)	97.0 (20.00-2.40) 95.8 (24.99-2.30)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.40 (at 2.31Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.210 , 0.280 0.205 , 0.270	Depositor DCC
R_{free} test set	1600 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	32.7	Xtriage
Anisotropy	0.636	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 45.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.024 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4880	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.16% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: TMC, 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	0.62	0/2381	1.30	15/3246 (0.5%)
1	B	0.62	0/2410	1.28	23/3286 (0.7%)
All	All	0.62	0/4791	1.29	38/6532 (0.6%)

There are no bond length outliers.

All (38) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	320	ARG	NE-CZ-NH2	-15.97	112.32	120.30
1	A	320	ARG	NE-CZ-NH1	12.53	126.56	120.30
1	A	237	ARG	CD-NE-CZ	9.97	137.55	123.60
1	B	237	ARG	CD-NE-CZ	8.59	135.62	123.60
1	B	237	ARG	NE-CZ-NH1	8.33	124.47	120.30
1	B	370	ARG	CD-NE-CZ	7.64	134.30	123.60
1	A	370	ARG	CD-NE-CZ	7.64	134.29	123.60
1	B	318	ARG	NE-CZ-NH1	7.31	123.95	120.30
1	A	46	MET	CA-CB-CG	6.92	125.06	113.30
1	B	130	MET	CA-CB-CG	6.80	124.85	113.30
1	B	132	TYR	CB-CG-CD1	6.79	125.07	121.00
1	A	293	ARG	CD-NE-CZ	6.64	132.90	123.60
1	B	87	TYR	CB-CG-CD2	6.61	124.97	121.00
1	A	176	ARG	NE-CZ-NH1	6.47	123.54	120.30
1	B	87	TYR	CB-CG-CD1	-6.44	117.14	121.00
1	A	222	ARG	CD-NE-CZ	6.38	132.53	123.60
1	B	236	ARG	NE-CZ-NH1	5.98	123.29	120.30
1	A	293	ARG	NE-CZ-NH2	-5.95	117.32	120.30
1	B	290	THR	N-CA-CB	5.88	121.48	110.30
1	B	373	GLY	N-CA-C	-5.88	98.41	113.10
1	B	163	ARG	NE-CZ-NH2	-5.70	117.45	120.30
1	B	80	TYR	CB-CG-CD2	5.64	124.39	121.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	80	TYR	CB-CG-CD1	-5.62	117.63	121.00
1	A	51	ARG	CD-NE-CZ	5.53	131.34	123.60
1	B	130	MET	CB-CG-SD	5.51	128.93	112.40
1	B	363	ASP	CB-CG-OD1	5.42	123.18	118.30
1	B	226	ARG	NE-CZ-NH2	5.37	122.98	120.30
1	A	296	GLU	OE1-CD-OE2	-5.32	116.91	123.30
1	A	307	VAL	CB-CA-C	-5.29	101.35	111.40
1	A	226	ARG	CD-NE-CZ	5.28	130.99	123.60
1	B	307	VAL	CG1-CB-CG2	5.27	119.33	110.90
1	B	363	ASP	CB-CG-OD2	-5.20	113.62	118.30
1	B	236	ARG	CD-NE-CZ	5.19	130.87	123.60
1	B	281	ARG	NE-CZ-NH2	-5.16	117.72	120.30
1	A	328	ASP	CB-CG-OD1	5.07	122.86	118.30
1	B	290	THR	CB-CA-C	-5.05	97.97	111.60
1	B	307	VAL	CB-CA-C	-5.04	101.83	111.40
1	A	172	TYR	CA-CB-CG	-5.03	103.85	113.40

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	2327	0	2347	72	0
1	B	2357	0	2374	66	0
2	A	5	0	0	1	0
2	B	5	0	0	0	0
3	A	18	0	16	1	0
3	B	18	0	16	1	0
4	A	77	0	0	7	0
4	B	73	0	0	7	0
All	All	4880	0	4753	136	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 14.

All (136) 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:126:ILE:HD11	1:B:126:ILE:HD11	1.46	0.95
1:A:130:MET:HB3	1:A:131:PRO:HD3	1.54	0.87
1:B:354:THR:HB	1:B:355:PRO:HD2	1.57	0.86
1:B:262:LEU:HD22	1:B:290:THR:HG22	1.60	0.81
1:A:262:LEU:HD22	1:A:290:THR:HG22	1.63	0.79
1:A:290:THR:HG21	4:A:2066:HOH:O	1.81	0.78
1:B:290:THR:HG21	4:B:2053:HOH:O	1.85	0.75
1:A:67:GLN:HG3	4:A:2008:HOH:O	1.86	0.74
1:A:153:PRO:HB2	1:A:154:PRO:HD3	1.70	0.72
1:B:66:THR:HG21	1:B:80:TYR:CE1	2.25	0.72
1:B:262:LEU:O	1:B:263:SER:HB2	1.88	0.71
1:A:285:GLY:HA2	1:A:290:THR:HG23	1.71	0.70
1:B:59:GLY:O	1:B:216:ARG:HD2	1.89	0.70
1:A:373:GLY:O	1:A:374:GLU:HB2	1.91	0.69
1:A:62:LYS:HG3	4:A:2004:HOH:O	1.91	0.69
1:B:140:ALA:HB3	1:B:141:PRO:HD3	1.76	0.69
1:B:358:ILE:HB	1:B:359:PRO:HD3	1.75	0.68
1:A:208:LEU:HD12	1:A:209:PRO:HD2	1.74	0.68
1:B:116:ASP:O	1:B:119:VAL:HG23	1.95	0.67
1:B:285:GLY:HA2	1:B:290:THR:HG23	1.77	0.66
1:A:212:ARG:NH1	1:A:216:ARG:HH21	1.94	0.66
1:A:354:THR:HB	1:A:355:PRO:HD2	1.78	0.66
1:A:153:PRO:CB	1:A:154:PRO:HD3	2.27	0.65
1:A:69:LEU:HD21	1:A:341:LEU:HD13	1.78	0.65
1:A:157:LEU:HD22	1:A:358:ILE:HG23	1.79	0.64
1:A:61:GLY:HA2	1:A:220:ARG:HH12	1.63	0.64
1:B:130:MET:HB3	1:B:131:PRO:HD3	1.80	0.64
1:A:285:GLY:HA2	1:A:290:THR:CG2	2.27	0.63
1:B:263:SER:HA	1:B:293:ARG:NH1	2.13	0.63
1:A:61:GLY:HA2	1:A:220:ARG:NH1	2.13	0.62
1:A:262:LEU:HD22	1:A:290:THR:CG2	2.32	0.59
1:B:285:GLY:HA2	1:B:290:THR:CG2	2.33	0.59
1:B:219:LYS:HA	4:B:2038:HOH:O	2.03	0.57
1:A:121:MET:HB3	1:A:182:MET:HE2	1.87	0.57
1:A:64:THR:O	1:A:68:LEU:HG	2.05	0.57
1:A:246:VAL:O	1:A:250:GLN:HG3	2.06	0.56
1:A:164:HIS:ND1	1:A:166:ILE:HG12	2.20	0.56
1:A:153:PRO:HB2	1:A:154:PRO:CD	2.36	0.56
1:A:69:LEU:HD11	1:A:341:LEU:HD13	1.87	0.55
1:B:333:PRO:HA	4:B:2005:HOH:O	2.06	0.55
1:A:364:LEU:HD22	1:B:310:TRP:CZ2	2.42	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:62:LYS:HB2	1:A:62:LYS:NZ	2.23	0.55
1:A:219:LYS:HB3	1:A:219:LYS:NZ	2.22	0.54
1:A:362:CYS:HB3	1:A:366:ARG:CZ	2.39	0.53
1:B:262:LEU:HD22	1:B:290:THR:CG2	2.36	0.52
1:B:124:ALA:O	1:B:128:MET:HG2	2.09	0.52
1:B:357:SER:O	1:B:361:ILE:HG13	2.10	0.52
1:A:299:ALA:HB1	1:A:300:PRO:HD2	1.91	0.52
1:B:109:GLN:HB3	1:B:111:GLU:HG3	1.92	0.51
1:B:86:THR:HB	1:B:372:MET:HA	1.91	0.51
1:A:221:GLN:OE1	1:A:225:GLU:HG2	2.11	0.51
1:B:164:HIS:CG	1:B:165:PRO:HD2	2.44	0.51
1:A:362:CYS:HB3	1:A:366:ARG:NH1	2.26	0.51
1:B:246:VAL:HG21	1:B:324:VAL:HG21	1.93	0.51
1:B:56:GLY:HA2	1:B:239:TYR:CD1	2.46	0.51
1:A:51:ARG:HG3	4:A:2001:HOH:O	2.11	0.50
1:B:66:THR:O	1:B:70:VAL:HG23	2.10	0.50
1:A:87:TYR:HA	1:A:91:LEU:HB2	1.94	0.50
1:A:130:MET:HB3	1:A:131:PRO:CD	2.33	0.50
1:A:159:LEU:HD13	1:A:161:PHE:CZ	2.46	0.49
1:A:216:ARG:HD3	4:A:2048:HOH:O	2.12	0.49
1:A:69:LEU:HD11	1:A:341:LEU:CD1	2.43	0.49
1:B:85:MET:CE	1:B:89:ARG:HG3	2.41	0.49
1:B:276:SER:HB2	1:B:324:VAL:HB	1.95	0.49
1:B:209:PRO:HD3	4:B:2063:HOH:O	2.12	0.49
1:B:240:GLY:HA2	1:B:274:PRO:HG2	1.95	0.48
1:B:59:GLY:HA3	1:B:217:LEU:HD13	1.95	0.48
1:B:367:THR:HA	1:B:370:ARG:NH1	2.29	0.48
1:B:88:TRP:HZ2	1:B:132:TYR:HH	1.61	0.48
1:B:373:GLY:O	1:B:374:GLU:C	2.52	0.48
1:A:327:LEU:HD23	1:A:329:TYR:CZ	2.48	0.48
1:B:217:LEU:CD2	1:B:232:LEU:HD13	2.44	0.48
1:B:97:ILE:HG23	1:B:101:PHE:HE2	1.79	0.48
1:A:212:ARG:HH12	1:A:216:ARG:HH21	1.62	0.48
1:B:354:THR:HB	1:B:355:PRO:CD	2.35	0.47
1:A:97:ILE:HD13	3:A:500:TMC:H152	1.95	0.47
1:A:58:HIS:HB3	1:A:163:ARG:NH2	2.30	0.47
1:B:331:GLN:HB2	4:B:2065:HOH:O	2.14	0.47
1:B:287:THR:HG1	1:B:289:PHE:HD2	1.60	0.47
1:A:79:VAL:HG12	1:A:80:TYR:N	2.29	0.47
1:A:69:LEU:CD2	1:A:341:LEU:HD13	2.45	0.47
1:B:62:LYS:NZ	1:B:62:LYS:HB2	2.29	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:171:CYS:HA	1:B:238:VAL:HG11	1.98	0.46
1:B:299:ALA:HA	1:B:305:TYR:CZ	2.51	0.46
1:A:213:HIS:NE2	1:A:232:LEU:HD11	2.31	0.46
1:B:247:ARG:NH1	1:B:275:GLN:O	2.49	0.46
1:B:46:MET:HA	1:B:46:MET:HE2	1.98	0.46
1:A:58:HIS:ND1	1:A:172:TYR:OH	2.43	0.45
1:A:178:LEU:HD21	1:A:291:LEU:HD22	1.97	0.45
1:B:366:ARG:HD2	4:B:2069:HOH:O	2.16	0.45
1:B:83:GLU:HG3	1:B:88:TRP:CH2	2.52	0.45
1:B:97:ILE:HD13	3:B:500:TMC:H152	1.98	0.45
1:A:139:LEU:HD21	1:A:159:LEU:HD11	1.99	0.45
1:A:358:ILE:N	1:A:359:PRO:HD2	2.32	0.45
1:B:74:SER:HB3	1:B:77:ASP:CG	2.37	0.45
1:B:182:MET:HG3	4:B:2018:HOH:O	2.17	0.44
1:A:130:MET:CB	1:A:131:PRO:HD3	2.35	0.44
1:A:222:ARG:O	1:A:225:GLU:HB3	2.17	0.44
1:B:69:LEU:HD21	1:B:341:LEU:CD1	2.47	0.44
1:B:102:THR:HA	1:B:226:ARG:HH12	1.83	0.44
1:A:148:GLY:O	1:A:149:SER:C	2.56	0.43
1:A:164:HIS:CG	1:A:165:PRO:HD2	2.53	0.43
1:B:66:THR:HG23	1:B:160:ILE:HG21	2.00	0.43
1:B:247:ARG:HH21	1:B:280:PRO:HD2	1.83	0.43
1:B:213:HIS:CE1	1:B:232:LEU:HD11	2.53	0.43
1:A:78:ILE:HD11	1:A:160:ILE:CD1	2.48	0.43
1:A:352:VAL:CG2	1:A:357:SER:HB2	2.48	0.43
1:B:121:MET:HG3	1:B:181:SER:HB2	2.01	0.43
1:B:218:ALA:O	1:B:219:LYS:HB2	2.18	0.43
1:A:68:LEU:HB2	1:A:337:ARG:HG3	2.00	0.42
1:B:109:GLN:CB	1:B:111:GLU:HG3	2.49	0.42
1:A:352:VAL:HG21	1:A:357:SER:HB2	2.00	0.42
1:A:212:ARG:HD3	4:A:2045:HOH:O	2.20	0.42
1:A:211:ASP:HB2	4:A:2046:HOH:O	2.19	0.42
1:B:57:PRO:HB3	1:B:235:ILE:HG23	2.02	0.42
1:A:220:ARG:HG3	1:A:220:ARG:O	2.20	0.41
1:A:299:ALA:HB1	1:A:300:PRO:CD	2.50	0.41
1:A:118:ALA:HA	1:A:181:SER:O	2.21	0.41
1:A:219:LYS:HB3	1:A:219:LYS:HZ2	1.85	0.41
1:A:304:LEU:HD22	1:A:308:PHE:HB2	2.02	0.41
1:A:154:PRO:HA	1:A:155:PRO:HD3	1.85	0.41
1:A:57:PRO:O	1:A:60:MET:HE2	2.21	0.41
1:B:78:ILE:HA	1:B:158:THR:O	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:87:TYR:HB2	1:B:372:MET:SD	2.59	0.41
1:B:106:ARG:NH1	1:B:111:GLU:OE1	2.54	0.41
1:A:222:ARG:HG3	1:A:222:ARG:HH11	1.85	0.41
1:B:46:MET:HE1	1:B:47:PRO:HD3	2.02	0.41
1:B:74:SER:HB3	1:B:77:ASP:OD1	2.21	0.41
1:A:58:HIS:HB2	2:A:400:SO4:O4	2.20	0.41
1:A:78:ILE:HD11	1:A:160:ILE:HD12	2.02	0.41
1:B:281:ARG:HA	1:B:282:PRO:HD3	1.81	0.41
1:B:74:SER:HB3	1:B:77:ASP:OD2	2.21	0.41
1:A:217:LEU:HD22	1:A:232:LEU:HD13	2.02	0.40
1:A:102:THR:O	1:A:106:ARG:HG3	2.21	0.40
1:A:254:SER:HB3	1:A:257:GLU:HB3	2.02	0.40
1:B:72:LEU:HD11	1:B:337:ARG:HD2	2.04	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	296/331 (89%)	277 (94%)	17 (6%)	2 (1%)	22	32
1	B	302/331 (91%)	288 (95%)	12 (4%)	2 (1%)	22	32
All	All	598/662 (90%)	565 (94%)	29 (5%)	4 (1%)	22	32

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	263	SER
1	A	154	PRO
1	A	320	ARG
1	B	147	ALA

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	246/264 (93%)	231 (94%)	15 (6%)	18	30
1	B	249/264 (94%)	237 (95%)	12 (5%)	25	41
All	All	495/528 (94%)	468 (94%)	27 (6%)	21	35

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	127	THR
1	A	130	MET
1	A	134	VAL
1	A	159	LEU
1	A	161	PHE
1	A	212	ARG
1	A	215	ASP
1	A	219	LYS
1	A	222	ARG
1	A	293	ARG
1	A	304	LEU
1	A	307	VAL
1	A	343	LEU
1	A	366	ARG
1	A	374	GLU
1	B	67	GLN
1	B	75	ARG
1	B	119	VAL
1	B	123	SER
1	B	130	MET
1	B	212	ARG
1	B	247	ARG
1	B	263	SER
1	B	307	VAL
1	B	366	ARG
1	B	368	PHE
1	B	374	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	349	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

4 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	400	-	4,4,4	0.59	0	6,6,6	0.18	0
2	SO4	B	400	-	4,4,4	0.73	0	6,6,6	0.51	0
3	TMC	B	500	-	18,20,20	2.21	3 (16%)	18,32,32	2.23	5 (27%)
3	TMC	A	500	-	18,20,20	2.09	2 (11%)	18,32,32	2.38	7 (38%)

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	TMC	A	500	-	-	4/7/29/29	0/3/3/3
3	TMC	B	500	-	-	4/7/29/29	0/3/3/3

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	500	TMC	C2-N8	-6.59	1.38	1.46
3	A	500	TMC	C2-N8	-5.69	1.39	1.46
3	B	500	TMC	C2-C4	-5.36	1.35	1.51
3	A	500	TMC	C2-C4	-4.84	1.36	1.51
3	B	500	TMC	C11-N8	2.25	1.38	1.35

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	500	TMC	C4-C2-N8	6.03	122.99	111.11
3	A	500	TMC	C4-C2-N8	5.44	121.83	111.11
3	A	500	TMC	C3-C4-C2	4.92	122.85	112.34
3	B	500	TMC	C3-C4-C2	4.34	121.61	112.34
3	A	500	TMC	C2-C4-C6	3.29	120.97	111.53
3	B	500	TMC	O16-C12-C1	-3.08	105.36	112.30
3	A	500	TMC	C14-C10-C1	2.92	107.65	103.22
3	A	500	TMC	O16-C12-C1	-2.87	105.84	112.30
3	B	500	TMC	C2-C4-C6	2.54	118.83	111.53
3	A	500	TMC	C1-C15-C13	2.22	61.06	59.95
3	A	500	TMC	O7-C6-N5	2.11	124.77	121.01
3	B	500	TMC	C1-C13-C17	-2.10	106.75	108.39

There are no chirality outliers.

All (8) torsion outliers are listed below:

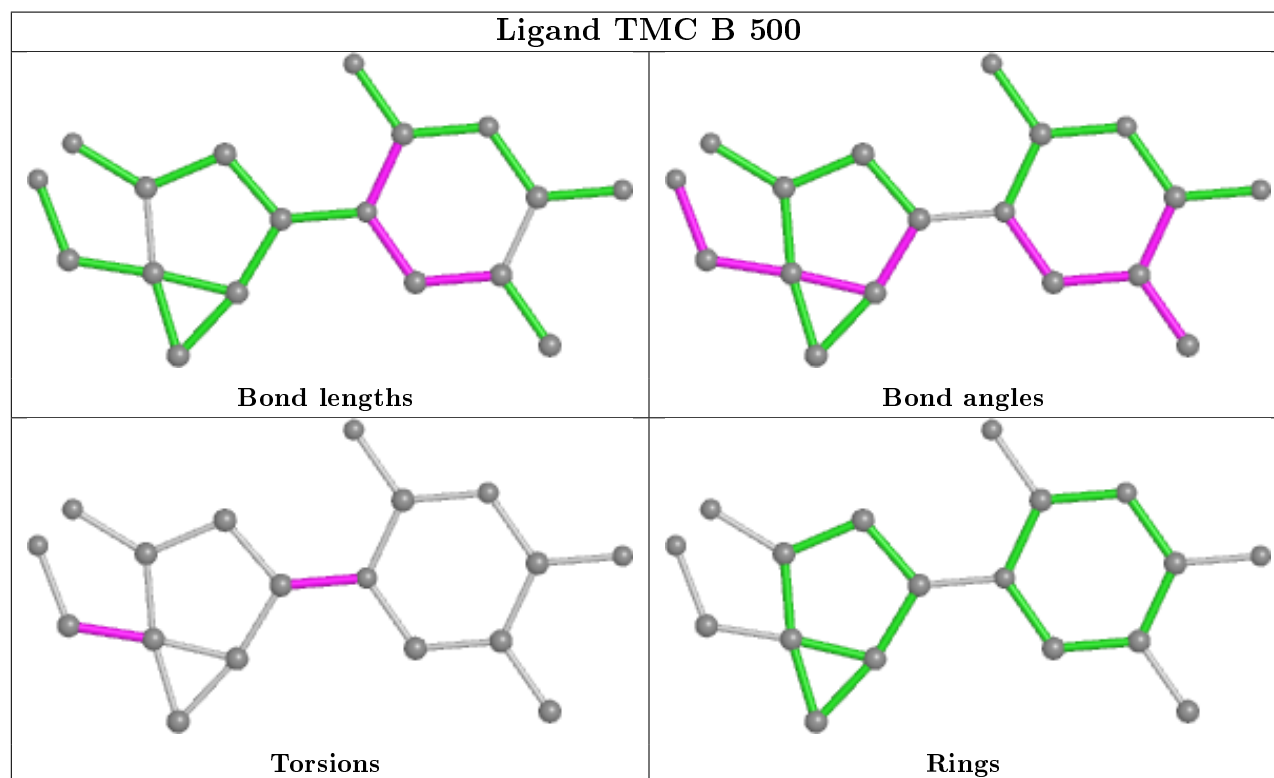
Mol	Chain	Res	Type	Atoms
3	B	500	TMC	C13-C17-N8-C2
3	B	500	TMC	C10-C1-C12-O16
3	A	500	TMC	C13-C17-N8-C2
3	A	500	TMC	C10-C1-C12-O16
3	B	500	TMC	C15-C1-C12-O16
3	A	500	TMC	C15-C1-C12-O16
3	B	500	TMC	C13-C1-C12-O16
3	A	500	TMC	C13-C1-C12-O16

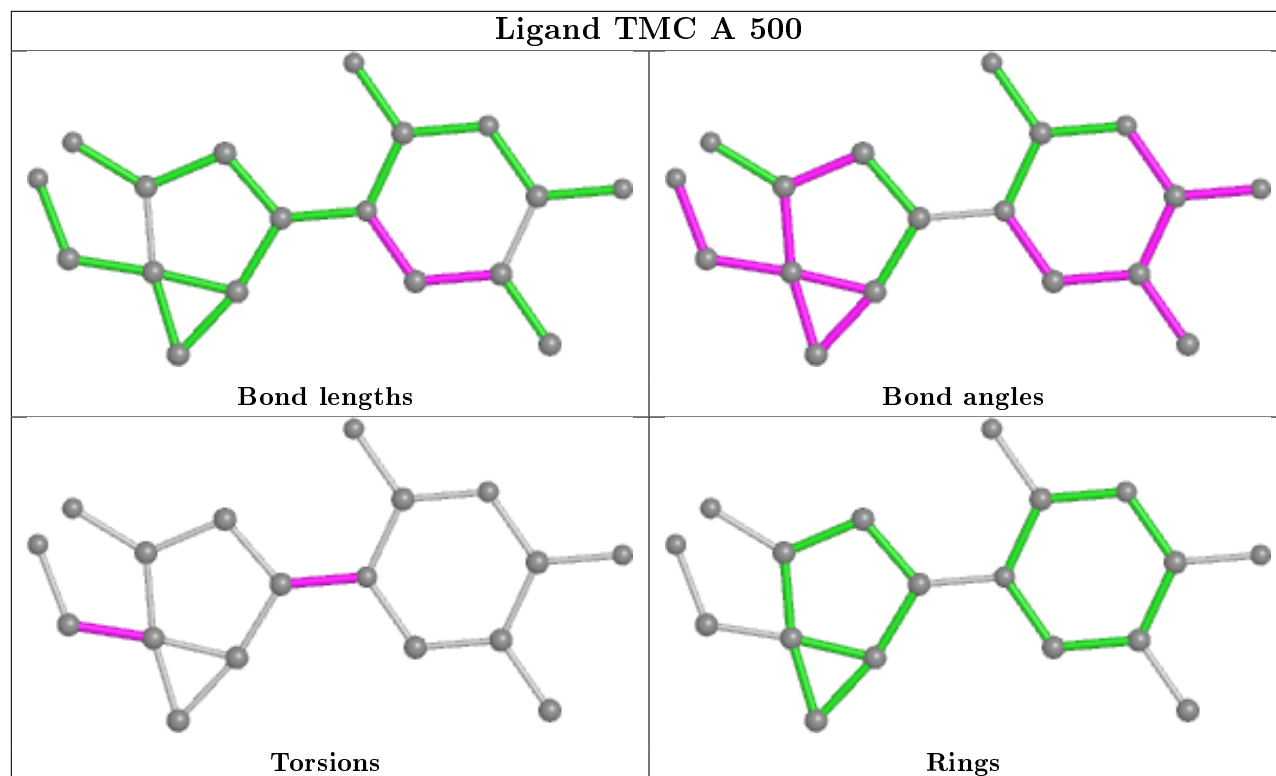
There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	400	SO4	1	0
3	B	500	TMC	1	0
3	A	500	TMC	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	304/331 (91%)	-0.20	13 (4%) 35 33	17, 32, 78, 110	0
1	B	310/331 (93%)	-0.21	17 (5%) 25 24	18, 33, 75, 138	0
All	All	614/662 (92%)	-0.21	30 (4%) 29 28	17, 33, 77, 138	0

All (30) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	226	ARG	5.8
1	B	274	PRO	5.5
1	B	264	GLY	5.4
1	A	148	GLY	5.2
1	A	149	SER	4.9
1	B	300	PRO	4.1
1	A	264	GLY	3.9
1	B	148	GLY	3.7
1	A	332	SER	3.7
1	A	68	LEU	3.6
1	A	65	THR	3.3
1	A	334	ALA	3.3
1	A	62	LYS	3.2
1	B	277	ASN	3.2
1	B	75	ARG	3.2
1	B	263	SER	3.0
1	A	263	SER	3.0
1	A	335	GLY	2.9
1	B	149	SER	2.9
1	B	217	LEU	2.8
1	B	46	MET	2.8
1	B	227	LEU	2.6
1	A	63	THR	2.6
1	B	355	PRO	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	276	SER	2.5
1	A	66	THR	2.4
1	B	275	GLN	2.4
1	B	215	ASP	2.4
1	A	69	LEU	2.2
1	B	74	SER	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

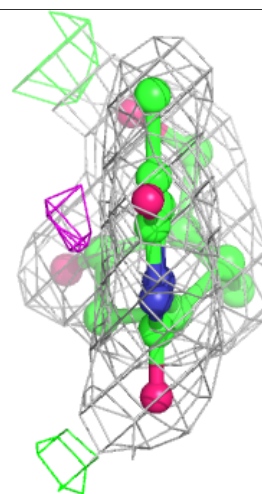
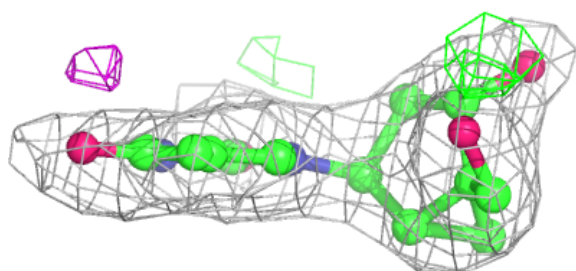
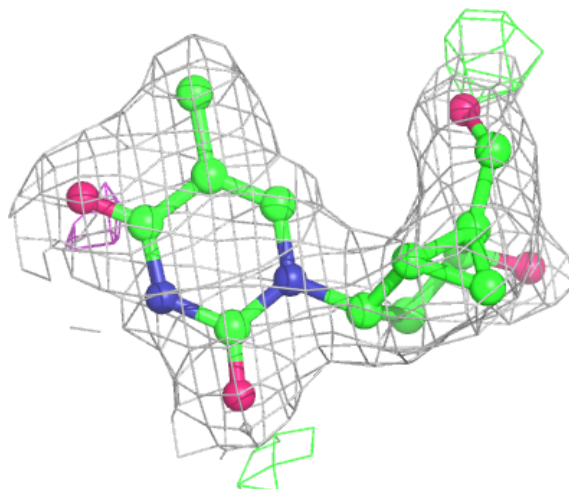
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

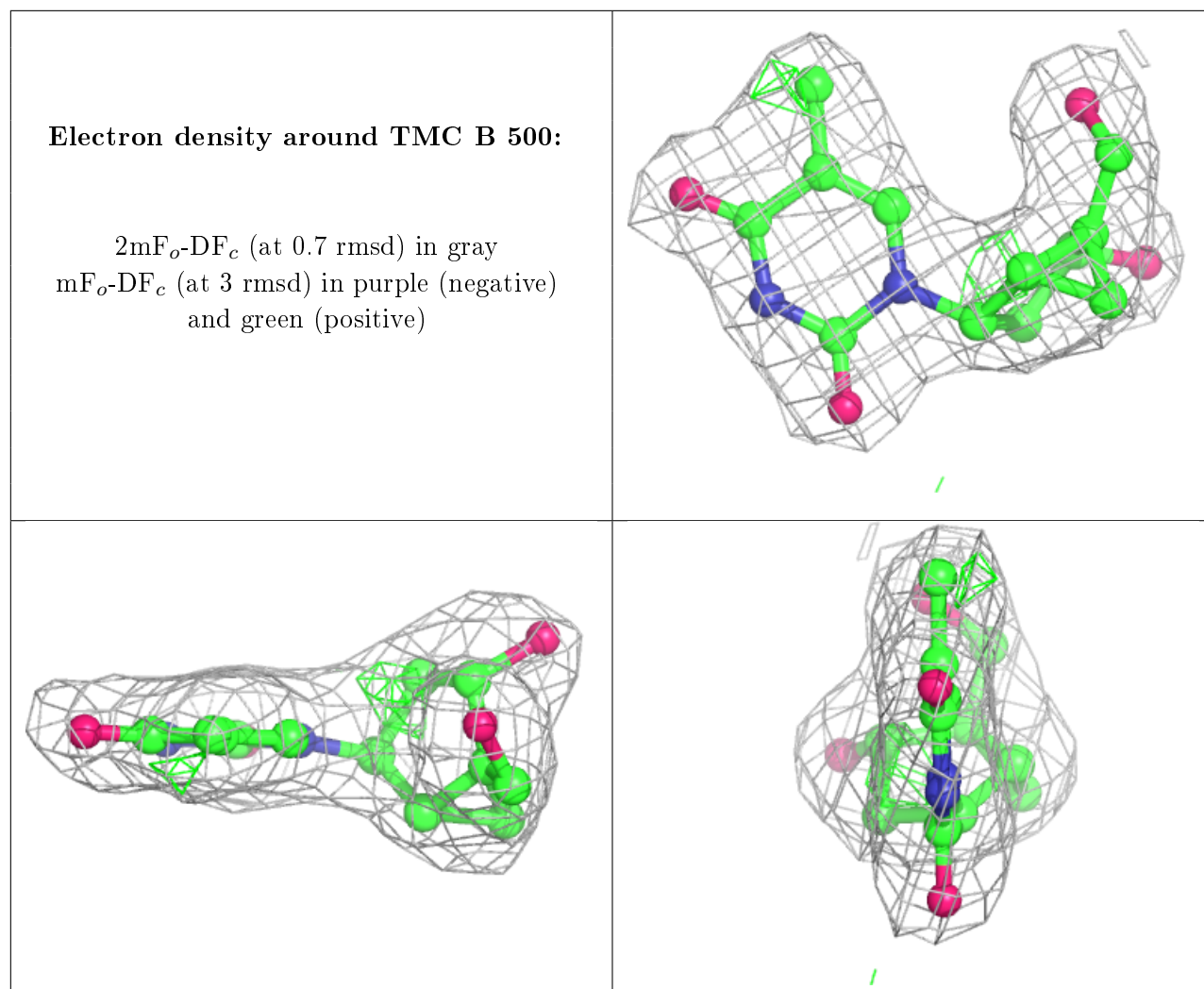
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	TMC	A	500	18/18	0.91	0.33	11,25,35,38	18
3	TMC	B	500	18/18	0.93	0.30	15,23,28,30	18
2	SO4	A	400	5/5	0.93	0.21	78,78,80,84	0
2	SO4	B	400	5/5	0.99	0.13	34,34,42,43	0

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 TMC A 500:

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





6.5 Other polymers ⓘ

There are no such residues in this entry.