



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 03:38 AM UTC

PDB ID : 8ICX / pdb_00008icx
Title : DNA POLYMERASE BETA (POL B) (E.C.2.7.7.7) COMPLEXED WITH SEVEN BASE PAIRS OF DNA; SOAKED IN THE PRESENCE OF DTTP (1 MILLIMOLAR) AND MNCL2 (5 MILLIMOLAR)
Authors : Pelletier, H.; Sawaya, M.R.
Deposited on : 1995-12-15
Resolution : 3.00 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)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

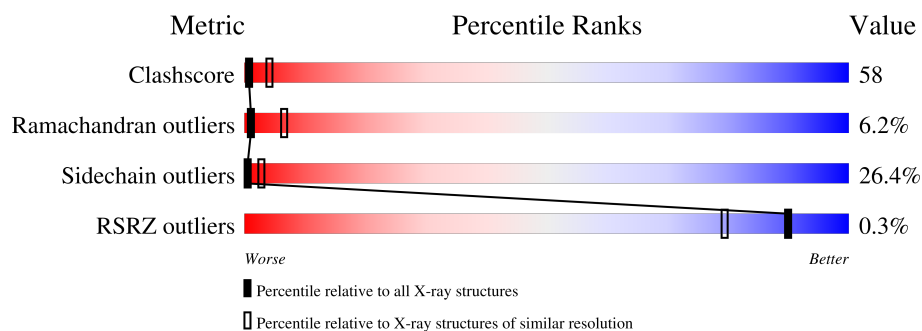
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)
RSRZ outliers	180081	2671 (3.00-3.00)

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 $\leq 5\%$. 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	T	8	 25% 38% 38%
2	P	7	 29% 71%
3	A	335	 21% 42% 27% 7% .

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 3070 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 DNA chain called DNA (5'-D(*CP*AP*TP*TP*AP*GP*AP*A)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	T	8	Total	C	N	O	P	0	0	0
			145	69	27	42	7			

- Molecule 2 is a DNA chain called DNA (5'-D(*TP*CP*TP*AP*AP*TP*G)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	P	7	Total	C	N	O	P	0	0	0
			144	69	24	44	7			

- Molecule 3 is a protein called PROTEIN (DNA POLYMERASE BETA (E.C.2.7.7.7)).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	A	327	Total	C	N	O	S	18	0	0
			2623	1657	458	499	9			

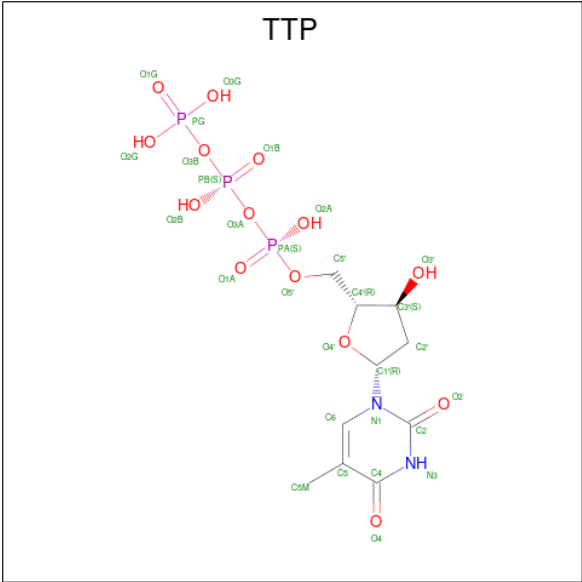
- Molecule 4 is MANGANESE (II) ION (CCD ID: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	2	Total	Mn	0	0
			2	2		

- Molecule 5 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	2	Total	Na	0	0
			2	2		

- Molecule 6 is THYMIDINE-5'-TRIPHOSPHATE (CCD ID: TTP) (formula: C₁₀H₁₇N₂O₁₄P₃).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	A	1	Total	C	O	P	0	0
			20	5	12	3		

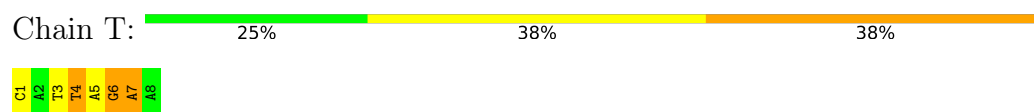
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	T	14	Total	O	0	0
			14	14		
7	P	17	Total	O	0	0
			17	17		
7	A	103	Total	O	0	0
			103	103		

3 Residue-property plots

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.

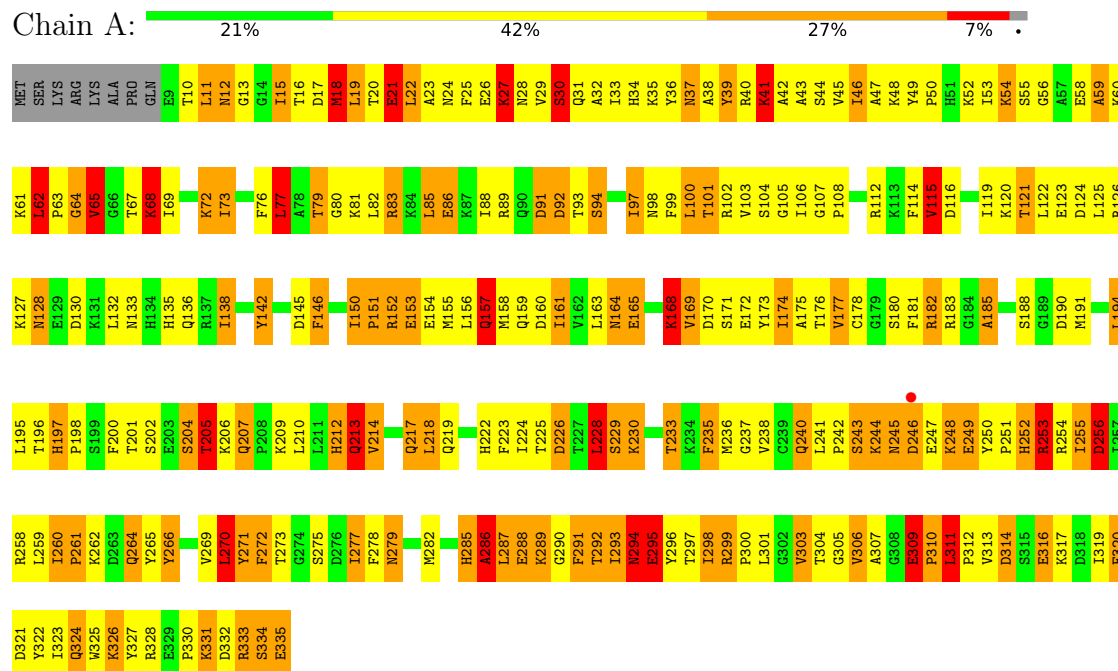
- Molecule 1: DNA (5'-D(*CP*AP*TP*TP*AP*GP*AP*A)-3')



- Molecule 2: DNA (5'-D(*TP*CP*TP*AP*AP*TP*G)-3')



- Molecule 3: PROTEIN (DNA POLYMERASE BETA (E.C.2.7.7.7))



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	179.69Å 57.67Å 48.13Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 3.00 20.00 – 3.00	Depositor EDS
% Data completeness (in resolution range)	91.0 (20.00-3.00) 90.5 (20.00-3.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.46 (at 2.70Å)	Xtriage
Refinement program	TNT 5-D	Depositor
R, R_{free}	0.161 , (Not available) (Not available) , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	36.5	Xtriage
Anisotropy	0.272	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.19 , 193.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	3070	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.92% 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 i

5.1 Standard geometry i

Bond lengths and bond angles in the following residue types are not validated in this section: NA, TTP, MN

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	T	1.03	0/162	3.66	12/249 (4.8%)
2	P	1.24	2/160 (1.2%)	3.83	15/243 (6.2%)
3	A	1.54	15/2672 (0.6%)	2.19	128/3590 (3.6%)
All	All	1.50	17/2994 (0.6%)	2.43	155/4082 (3.8%)

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
3	A	2	0

All (17) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	217	GLN	N-CA	-7.05	1.38	1.46
2	P	1	DT	OP3-P	6.32	1.61	1.48
3	A	170	ASP	CA-C	-6.15	1.45	1.52
3	A	242	PRO	N-CA	-6.15	1.39	1.47
3	A	242	PRO	CA-C	-5.95	1.46	1.52
3	A	260	ILE	CA-C	-5.93	1.46	1.52
3	A	170	ASP	N-CA	-5.83	1.39	1.46
3	A	63	PRO	N-CA	-5.78	1.40	1.47
3	A	165	GLU	N-CA	-5.69	1.39	1.46
3	A	24	ASN	CA-C	-5.32	1.45	1.52
3	A	224	ILE	CA-CB	-5.31	1.47	1.54
3	A	15	ILE	CA-C	-5.25	1.46	1.52
2	P	2	DC	C1'-N1	5.21	1.65	1.49
3	A	224	ILE	CA-C	-5.21	1.46	1.52
3	A	291	PHE	CA-C	-5.16	1.46	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	261	PRO	CA-C	-5.13	1.47	1.52
3	A	326	LYS	CE-NZ	-5.11	1.34	1.49

All (155) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	T	7	DA	C4-N9-C1'	-25.21	89.23	127.05
1	T	4	DT	C6-N1-C1'	-23.99	83.36	119.35
1	T	4	DT	C2-N1-C1'	23.36	154.40	119.35
1	T	7	DA	C8-N9-C1'	23.35	162.08	127.05
2	P	1	DT	C6-N1-C1'	-20.65	88.37	119.35
2	P	1	DT	C2-N1-C1'	20.22	149.68	119.35
2	P	2	DC	C2-N1-C1'	16.47	144.40	119.70
2	P	6	DT	C6-N1-C1'	-15.42	96.22	119.35
2	P	7	DG	C4-N9-C1'	-15.25	104.12	127.00
2	P	6	DT	C2-N1-C1'	14.73	141.44	119.35
2	P	2	DC	C6-N1-C1'	-14.68	97.68	119.70
2	P	7	DG	C8-N9-C1'	14.55	148.83	127.00
2	P	5	DA	C8-N9-C1'	-13.34	107.04	127.05
2	P	5	DA	C4-N9-C1'	12.26	145.44	127.05
2	P	3	DT	C2-N1-C1'	12.10	137.51	119.35
2	P	3	DT	C6-N1-C1'	-11.42	102.23	119.35
3	A	49	TYR	CA-C-N	11.23	133.88	119.84
3	A	49	TYR	C-N-CA	11.23	133.88	119.84
1	T	6	DG	C4-N9-C1'	-10.23	111.66	127.00
3	A	49	TYR	O-C-N	9.94	129.23	121.47
3	A	256	ASP	CA-CB-CG	9.83	122.43	112.60
1	T	6	DG	C8-N9-C1'	9.13	140.70	127.00
3	A	83	ARG	N-CA-CB	8.86	122.85	110.01
1	T	3	DT	P-O5'-C5'	8.82	133.23	120.00
3	A	256	ASP	N-CA-CB	8.60	125.26	110.81
3	A	68	LYS	N-CA-CB	8.57	122.85	109.82
3	A	40	ARG	N-CA-C	8.03	120.11	111.36
3	A	168	LYS	N-CA-CB	7.92	121.76	110.12
3	A	291	PHE	CA-CB-CG	-7.87	105.93	113.80
3	A	266	TYR	CA-CB-CG	-7.86	99.75	113.90
3	A	130	ASP	CA-CB-CG	7.68	120.28	112.60
2	P	3	DT	N1-C1'-C2'	7.67	125.00	113.50
3	A	150	ILE	N-CA-C	7.62	115.10	107.55
3	A	309	GLU	CA-C-N	7.48	129.19	119.84
3	A	309	GLU	C-N-CA	7.48	129.19	119.84
3	A	235	PHE	CA-CB-CG	-7.45	106.35	113.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	213	GLN	CB-CA-C	-7.43	98.92	110.81
3	A	88	ILE	CB-CA-C	-7.35	102.39	112.02
3	A	272	PHE	CA-CB-CG	-7.33	106.47	113.80
3	A	157	GLN	N-CA-CB	7.19	120.80	110.16
3	A	314	ASP	N-CA-CB	7.12	121.14	110.53
3	A	311	LEU	CA-C-N	7.11	128.98	120.23
3	A	311	LEU	C-N-CA	7.11	128.98	120.23
3	A	20	THR	N-CA-CB	7.10	120.37	110.07
3	A	65	VAL	CB-CA-C	7.08	121.52	111.81
3	A	20	THR	CB-CA-C	7.08	122.08	110.90
3	A	286	ALA	CA-C-N	7.01	129.97	120.44
3	A	286	ALA	C-N-CA	7.01	129.97	120.44
3	A	214	VAL	N-CA-CB	6.83	119.83	110.54
3	A	32	ALA	N-CA-C	6.81	121.25	110.70
3	A	142	TYR	CB-CA-C	-6.80	100.52	111.06
3	A	333	ARG	N-CA-C	-6.80	104.86	113.02
3	A	30	SER	CA-C-N	-6.77	112.41	122.83
3	A	30	SER	C-N-CA	-6.77	112.41	122.83
3	A	45	VAL	N-CA-C	6.67	116.80	110.53
3	A	228	LEU	N-CA-C	-6.66	102.45	112.04
1	T	1	DC	P-O3'-C3'	6.62	130.14	120.20
3	A	12	ASN	CB-CA-C	6.62	121.15	110.09
3	A	25	PHE	CA-CB-CG	-6.62	107.19	113.80
3	A	229	SER	CA-C-O	6.57	128.00	119.98
1	T	7	DA	N9-C1'-C2'	-6.49	103.76	113.50
3	A	252	HIS	CA-CB-CG	-6.46	107.34	113.80
3	A	313	VAL	N-CA-C	6.45	117.15	107.80
3	A	291	PHE	N-CA-C	6.43	116.76	108.34
3	A	98	ASN	N-CA-C	6.39	118.24	111.28
3	A	320	PHE	CA-CB-CG	-6.37	107.43	113.80
3	A	27	LYS	N-CA-C	6.30	118.95	111.71
3	A	102	ARG	CA-CB-CG	-6.25	101.60	114.10
2	P	3	DT	O4'-C1'-N1	6.24	117.76	108.40
3	A	62	LEU	N-CA-C	6.23	117.65	109.93
3	A	312	PRO	CB-CA-C	-6.17	104.83	111.56
3	A	159	GLN	CB-CA-C	-6.16	100.57	110.79
3	A	65	VAL	N-CA-C	6.10	116.16	106.88
3	A	65	VAL	N-CA-CB	6.10	118.92	111.60
3	A	142	TYR	CA-CB-CG	-6.09	102.93	113.90
3	A	303	VAL	N-CA-C	6.07	120.63	111.89
3	A	201	THR	N-CA-CB	-6.07	101.78	111.43
3	A	116	ASP	CA-CB-CG	-5.92	106.68	112.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	270	LEU	O-C-N	5.91	128.40	122.08
3	A	116	ASP	N-CA-CB	5.89	118.56	110.01
3	A	73	ILE	N-CA-CB	5.87	117.03	110.51
3	A	294	ASN	CA-CB-CG	5.87	118.47	112.60
3	A	170	ASP	N-CA-C	-5.85	98.27	107.93
3	A	243	SER	CB-CA-C	5.85	119.56	110.67
3	A	292	THR	CB-CA-C	5.84	119.73	109.80
3	A	161	ILE	N-CA-C	5.84	116.57	110.62
3	A	41	LYS	CB-CA-C	5.84	120.77	110.85
3	A	12	ASN	N-CA-CB	5.80	119.07	110.49
3	A	309	GLU	N-CA-C	5.78	122.59	109.81
3	A	229	SER	CA-C-N	-5.77	114.63	122.77
3	A	229	SER	C-N-CA	-5.77	114.63	122.77
3	A	91	ASP	N-CA-CB	5.72	120.16	110.49
3	A	77	LEU	N-CA-CB	5.71	118.61	110.16
3	A	68	LYS	O-C-N	5.70	128.18	122.08
3	A	59	ALA	O-C-N	5.69	127.93	122.07
3	A	219	GLN	CB-CA-C	-5.69	98.78	110.38
3	A	288	GLU	N-CA-C	-5.68	105.01	111.14
3	A	27	LYS	CB-CA-C	5.67	121.13	110.63
3	A	151	PRO	N-CA-CB	5.67	108.37	103.27
3	A	243	SER	N-CA-C	-5.66	101.00	109.15
3	A	224	ILE	CA-CB-CG1	-5.64	100.81	110.40
3	A	271	TYR	CA-CB-CG	-5.62	103.78	113.90
3	A	115	VAL	O-C-N	5.61	127.40	121.91
3	A	177	VAL	N-CA-C	-5.59	99.68	108.23
3	A	21	GLU	CA-C-N	5.58	128.55	120.79
3	A	21	GLU	C-N-CA	5.58	128.55	120.79
3	A	197	HIS	O-C-N	5.56	125.71	121.65
3	A	24	ASN	N-CA-CB	5.53	118.88	110.14
3	A	334	SER	N-CA-C	5.52	118.81	111.75
3	A	261	PRO	N-CA-C	-5.46	102.11	110.95
3	A	165	GLU	O-C-N	5.45	127.98	122.09
3	A	223	PHE	N-CA-C	-5.43	104.90	112.45
3	A	11	LEU	N-CA-C	-5.43	105.79	112.90
3	A	35	LYS	CA-C-O	5.43	126.49	120.63
3	A	190	ASP	CA-C-N	-5.43	114.54	122.19
3	A	190	ASP	C-N-CA	-5.43	114.54	122.19
3	A	253	ARG	CA-C-O	5.42	126.56	120.70
3	A	212	HIS	CA-CB-CG	-5.39	108.41	113.80
3	A	17	ASP	N-CA-CB	5.38	118.13	110.16
3	A	64	GLY	CA-C-N	-5.38	116.27	122.95

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	64	GLY	C-N-CA	-5.38	116.27	122.95
3	A	101	THR	CB-CA-C	5.38	120.58	110.63
3	A	246	ASP	N-CA-C	5.38	122.25	110.80
3	A	49	TYR	CB-CA-C	5.36	117.70	109.60
2	P	7	DG	P-O5'-C5'	-5.33	112.00	120.00
3	A	17	ASP	CB-CA-C	5.30	119.86	110.85
3	A	18	MET	N-CA-C	-5.27	105.62	111.36
1	T	4	DT	O4'-C1'-C2'	-5.26	98.52	106.40
3	A	165	GLU	CA-C-O	-5.25	115.29	120.70
3	A	11	LEU	CA-C-N	-5.23	113.91	122.54
3	A	11	LEU	C-N-CA	-5.23	113.91	122.54
1	T	3	DT	O4'-C1'-N1	5.22	116.24	108.40
3	A	253	ARG	N-CA-C	5.21	117.89	109.40
3	A	261	PRO	O-C-N	5.21	128.66	122.98
3	A	68	LYS	CB-CA-C	5.20	119.19	110.92
3	A	229	SER	O-C-N	-5.20	116.81	123.26
3	A	285	HIS	N-CA-C	5.17	116.78	111.03
3	A	115	VAL	N-CA-CB	-5.16	104.51	110.55
3	A	54	LYS	N-CA-C	5.16	120.22	113.30
3	A	98	ASN	CB-CA-C	5.16	119.35	110.79
3	A	176	THR	CA-C-N	-5.15	115.93	122.37
3	A	176	THR	C-N-CA	-5.15	115.93	122.37
1	T	4	DT	C1'-O4'-C4'	-5.13	102.00	109.70
3	A	62	LEU	CA-C-O	5.10	125.43	120.23
3	A	299	ARG	CB-CA-C	5.10	115.77	108.68
3	A	240	GLN	CB-CG-CD	-5.10	103.94	112.60
3	A	86	GLU	N-CA-CB	5.09	117.40	110.01
3	A	291	PHE	CB-CA-C	-5.09	101.12	113.19
3	A	64	GLY	N-CA-C	-5.08	108.31	114.92
3	A	39	TYR	CA-CB-CG	-5.03	104.84	113.90
3	A	146	PHE	CA-CB-CG	-5.02	108.78	113.80
3	A	164	ASN	CA-CB-CG	-5.01	107.58	112.60
3	A	170	ASP	O-C-N	5.01	129.47	123.26
3	A	233	THR	CA-CB-OG1	-5.01	102.08	109.60
3	A	295	GLU	CB-CG-CD	5.01	121.11	112.60

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	A	68	LYS	CA
3	A	246	ASP	CA

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	T	145	0	80	5	0
2	P	144	0	81	17	0
3	A	2623	0	2641	308	0
4	A	2	0	0	0	0
5	A	2	0	0	0	0
6	A	20	0	7	5	0
7	A	103	0	0	19	0
7	P	17	0	0	1	0
7	T	14	0	0	3	0
All	All	3070	0	2809	329	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 58.

All (329) 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)
3:A:293:ILE:HD13	3:A:298:ILE:HG13	1.30	1.11
3:A:180:SER:HB3	3:A:183:ARG:HH21	1.18	1.07
3:A:245:ASN:N	3:A:245:ASN:HD22	1.55	1.04
2:P:1:DT:H2''	2:P:2:DC:H5'	1.41	1.01
3:A:277:ILE:HD13	3:A:277:ILE:H	1.26	1.01
3:A:60:LYS:HA	3:A:65:VAL:HG23	1.47	0.94
3:A:243:SER:HB3	3:A:249:GLU:HA	1.52	0.90
3:A:151:PRO:HG2	3:A:154:GLU:HG3	1.52	0.88
3:A:18:MET:HG2	3:A:19:LEU:N	1.89	0.88
3:A:150:ILE:HG21	3:A:158:MET:HE1	1.57	0.87
1:T:6:DG:H2''	1:T:7:DA:C8	2.10	0.87
3:A:29:VAL:HG21	3:A:94:SER:HB2	1.57	0.86
3:A:150:ILE:CD1	3:A:253:ARG:HG2	2.06	0.85
2:P:5:DA:H2''	2:P:6:DT:H5''	1.59	0.84
3:A:197:HIS:CD2	3:A:198:PRO:HD2	2.11	0.84

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:41:LYS:HE2	3:A:64:GLY:HA3	1.60	0.83
3:A:18:MET:HE1	3:A:76:PHE:HB2	1.60	0.83
3:A:182:ARG:HG2	3:A:182:ARG:HH11	1.44	0.82
3:A:68:LYS:HB2	3:A:68:LYS:NZ	1.92	0.82
2:P:5:DA:H2''	2:P:6:DT:C5'	2.10	0.81
3:A:18:MET:HE2	3:A:82:LEU:HD22	1.62	0.80
3:A:29:VAL:HA	3:A:97:ILE:HD12	1.63	0.80
3:A:243:SER:CB	3:A:249:GLU:HA	2.13	0.79
3:A:182:ARG:HG2	3:A:182:ARG:NH1	1.97	0.79
3:A:121:THR:HG23	3:A:124:ASP:CG	2.07	0.79
3:A:277:ILE:HG12	3:A:335:GLU:HA	1.64	0.79
3:A:285:HIS:NE2	3:A:289:LYS:HG2	1.98	0.78
3:A:60:LYS:HA	3:A:65:VAL:CG2	2.13	0.77
3:A:125:LEU:CD2	3:A:132:LEU:HD21	2.16	0.76
3:A:23:ALA:HB2	3:A:39:TYR:HB3	1.66	0.76
3:A:178:CYS:SG	3:A:194:LEU:HD22	2.26	0.76
3:A:73:ILE:HG22	3:A:77:LEU:CD2	2.15	0.75
3:A:270:LEU:HD23	3:A:319:ILE:HG21	1.65	0.75
3:A:165:GLU:HA	3:A:168:LYS:HG2	1.68	0.75
3:A:306:VAL:HG22	7:A:650:HOH:O	1.85	0.74
3:A:286:ALA:CB	3:A:323:ILE:HG21	2.18	0.73
3:A:210:LEU:CB	3:A:259:LEU:HD21	2.17	0.73
3:A:245:ASN:N	3:A:245:ASN:ND2	2.29	0.73
3:A:272:PHE:HD1	6:A:338:TTP:H4'	1.54	0.73
3:A:21:GLU:OE1	3:A:85:LEU:HG	1.89	0.72
3:A:286:ALA:CB	3:A:293:ILE:HD11	2.20	0.72
3:A:180:SER:HB3	3:A:183:ARG:NH2	1.98	0.72
3:A:270:LEU:HD13	3:A:316:GLU:OE1	1.89	0.72
3:A:180:SER:CB	3:A:183:ARG:HH21	1.99	0.72
3:A:37:ASN:HB3	7:A:556:HOH:O	1.90	0.71
3:A:323:ILE:O	3:A:324:GLN:HG2	1.90	0.71
3:A:11:LEU:HD23	3:A:11:LEU:H	1.53	0.71
3:A:300:PRO:HD3	3:A:311:LEU:HD22	1.73	0.70
3:A:210:LEU:HB3	3:A:259:LEU:HD21	1.72	0.70
3:A:41:LYS:HE2	3:A:64:GLY:CA	2.21	0.70
3:A:127:LYS:HB2	3:A:128:ASN:ND2	2.07	0.70
3:A:155:MET:HE2	3:A:188:SER:CB	2.22	0.70
3:A:27:LYS:HB3	3:A:36:TYR:CG	2.27	0.69
3:A:260:ILE:HG23	3:A:261:PRO:HD2	1.74	0.69
3:A:12:ASN:HB3	3:A:46:ILE:HD12	1.73	0.69
3:A:112:ARG:O	3:A:115:VAL:HB	1.91	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:108:PRO:O	3:A:112:ARG:HG3	1.93	0.69
3:A:128:ASN:ND2	3:A:128:ASN:N	2.41	0.69
3:A:165:GLU:HA	3:A:168:LYS:CG	2.21	0.69
3:A:92:ASP:HB3	7:A:647:HOH:O	1.91	0.69
3:A:121:THR:HG23	3:A:124:ASP:OD1	1.91	0.69
3:A:18:MET:HE1	3:A:76:PHE:CB	2.22	0.69
3:A:233:THR:HB	7:A:538:HOH:O	1.93	0.68
3:A:212:HIS:HB3	7:A:541:HOH:O	1.93	0.68
3:A:123:GLU:HG3	7:A:623:HOH:O	1.92	0.68
3:A:245:ASN:HD22	3:A:245:ASN:H	1.43	0.67
3:A:180:SER:HA	3:A:183:ARG:HE	1.58	0.67
3:A:197:HIS:O	3:A:262:LYS:HB2	1.95	0.67
3:A:23:ALA:HB2	3:A:39:TYR:CB	2.25	0.67
3:A:31:GLN:HB2	3:A:112:ARG:HH12	1.59	0.67
3:A:79:THR:O	3:A:81:LYS:N	2.28	0.66
3:A:294:ASN:O	3:A:296:TYR:N	2.27	0.66
3:A:73:ILE:HG22	3:A:77:LEU:HD21	1.76	0.66
3:A:286:ALA:HB1	3:A:293:ILE:HD11	1.78	0.66
3:A:286:ALA:HB2	3:A:323:ILE:HG21	1.78	0.66
3:A:327:TYR:HD1	3:A:328:ARG:N	1.94	0.66
3:A:15:ILE:CG2	3:A:46:ILE:HD13	2.25	0.65
3:A:128:ASN:N	3:A:128:ASN:HD22	1.95	0.65
3:A:155:MET:HA	3:A:158:MET:HE3	1.79	0.65
3:A:293:ILE:HD13	3:A:298:ILE:CG1	2.19	0.65
3:A:19:LEU:HB3	3:A:43:ALA:HB2	1.79	0.65
3:A:150:ILE:HG21	3:A:158:MET:CE	2.27	0.64
3:A:286:ALA:O	3:A:291:PHE:N	2.30	0.64
1:T:5:DA:H2''	1:T:6:DG:O5'	1.96	0.64
3:A:41:LYS:HD3	3:A:42:ALA:H	1.63	0.64
3:A:150:ILE:HD11	3:A:253:ARG:HG2	1.79	0.63
2:P:1:DT:H2''	2:P:2:DC:C5'	2.23	0.63
3:A:11:LEU:H	3:A:11:LEU:CD2	2.11	0.63
3:A:142:TYR:CE2	3:A:238:VAL:HG11	2.34	0.63
3:A:114:PHE:CZ	3:A:132:LEU:HD23	2.34	0.63
3:A:151:PRO:HG2	3:A:154:GLU:CG	2.24	0.63
3:A:183:ARG:HH11	3:A:275:SER:HA	1.63	0.62
3:A:197:HIS:CG	3:A:198:PRO:HD2	2.34	0.62
3:A:182:ARG:NH1	3:A:273:THR:HG21	2.15	0.62
2:P:6:DT:H2''	2:P:7:DG:H5''	1.82	0.62
3:A:12:ASN:HD21	3:A:53:ILE:H	1.45	0.62
3:A:326:LYS:HG3	3:A:326:LYS:O	2.00	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:41:LYS:NZ	3:A:64:GLY:O	2.32	0.62
3:A:73:ILE:HG22	3:A:77:LEU:HD22	1.80	0.62
3:A:260:ILE:CG2	3:A:261:PRO:HD2	2.30	0.61
2:P:5:DA:C2'	2:P:6:DT:H5''	2.30	0.61
3:A:41:LYS:HD3	3:A:42:ALA:N	2.15	0.61
3:A:27:LYS:HB3	3:A:36:TYR:CD1	2.35	0.61
3:A:240:GLN:NE2	3:A:250:TYR:O	2.28	0.61
3:A:41:LYS:O	3:A:44:SER:HB3	2.01	0.60
3:A:278:PHE:CE2	3:A:333:ARG:HD2	2.35	0.60
2:P:1:DT:C2'	2:P:2:DC:H5'	2.24	0.60
3:A:288:GLU:OE1	3:A:288:GLU:HA	2.01	0.60
3:A:323:ILE:C	3:A:324:GLN:HG2	2.26	0.60
3:A:123:GLU:O	3:A:127:LYS:HG2	2.01	0.60
3:A:213:GLN:HG3	7:A:644:HOH:O	2.00	0.60
3:A:27:LYS:HE3	3:A:28:ASN:HD21	1.67	0.60
3:A:146:PHE:HZ	3:A:238:VAL:HG22	1.65	0.60
3:A:154:GLU:O	3:A:158:MET:HG3	2.02	0.60
3:A:294:ASN:HB2	3:A:295:GLU:OE1	2.02	0.60
3:A:210:LEU:HB2	3:A:259:LEU:HD21	1.83	0.59
3:A:18:MET:CE	3:A:76:PHE:HB2	2.32	0.59
3:A:155:MET:HE2	3:A:188:SER:HB2	1.82	0.59
3:A:180:SER:HB2	3:A:185:ALA:HB2	1.83	0.59
3:A:145:ASP:HB3	3:A:252:HIS:O	2.02	0.59
3:A:15:ILE:HG21	3:A:46:ILE:HD13	1.84	0.59
3:A:237:GLY:O	3:A:254:ARG:NH1	2.35	0.58
3:A:119:ILE:CG2	3:A:124:ASP:HB2	2.33	0.58
3:A:200:PHE:O	3:A:262:LYS:N	2.32	0.58
3:A:292:THR:O	3:A:298:ILE:HA	2.04	0.58
3:A:27:LYS:HG2	3:A:36:TYR:CE1	2.39	0.58
3:A:119:ILE:N	3:A:119:ILE:HD13	2.19	0.58
3:A:259:LEU:O	3:A:260:ILE:HD13	2.04	0.58
3:A:200:PHE:O	3:A:261:PRO:HA	2.03	0.58
3:A:150:ILE:HD13	3:A:253:ARG:HG2	1.85	0.58
3:A:182:ARG:NH2	3:A:316:GLU:OE2	2.34	0.58
3:A:209:LYS:HA	3:A:212:HIS:HB2	1.86	0.58
3:A:319:ILE:O	3:A:322:TYR:HB2	2.04	0.57
3:A:119:ILE:HG22	3:A:124:ASP:HB2	1.87	0.57
3:A:172:GLU:HB3	3:A:197:HIS:NE2	2.20	0.57
3:A:277:ILE:H	3:A:277:ILE:CD1	2.08	0.57
3:A:68:LYS:HB2	3:A:68:LYS:HZ1	1.70	0.56
3:A:103:VAL:HA	7:A:600:HOH:O	2.05	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:18:MET:HG3	3:A:22:LEU:HD23	1.88	0.56
3:A:181:PHE:HA	7:A:530:HOH:O	2.04	0.56
3:A:303:VAL:C	3:A:305:GLY:H	2.12	0.56
3:A:286:ALA:HA	3:A:323:ILE:CG2	2.36	0.55
1:T:6:DG:N7	7:T:652:HOH:O	2.33	0.55
3:A:18:MET:O	3:A:21:GLU:HB2	2.07	0.55
3:A:289:LYS:HD3	3:A:289:LYS:N	2.20	0.55
3:A:326:LYS:O	3:A:328:ARG:HG2	2.06	0.55
3:A:331:LYS:HG2	3:A:332:ASP:N	2.20	0.55
3:A:31:GLN:HE21	3:A:112:ARG:HH12	1.53	0.55
3:A:125:LEU:HD22	3:A:132:LEU:HD21	1.89	0.55
3:A:151:PRO:HB2	3:A:153:GLU:HG2	1.88	0.54
3:A:271:TYR:HB2	7:A:592:HOH:O	2.07	0.54
3:A:121:THR:OG1	3:A:122:LEU:N	2.40	0.54
3:A:154:GLU:HB3	3:A:241:LEU:HD11	1.91	0.53
3:A:121:THR:HG23	3:A:124:ASP:OD2	2.08	0.53
3:A:97:ILE:HD13	3:A:112:ARG:HG2	1.89	0.53
3:A:270:LEU:C	3:A:270:LEU:HD12	2.31	0.53
3:A:188:SER:HB3	7:A:522:HOH:O	2.08	0.53
3:A:83:ARG:HA	3:A:86:GLU:HG2	1.90	0.53
3:A:133:ASN:ND2	3:A:135:HIS:H	2.06	0.53
3:A:270:LEU:CD2	3:A:319:ILE:HG21	2.35	0.53
3:A:182:ARG:NH1	3:A:273:THR:OG1	2.42	0.52
3:A:244:LYS:HB3	3:A:245:ASN:ND2	2.24	0.52
3:A:330:PRO:O	3:A:333:ARG:HG2	2.09	0.52
3:A:295:GLU:HA	7:A:592:HOH:O	2.08	0.52
3:A:295:GLU:HG2	3:A:296:TYR:CE1	2.44	0.52
3:A:68:LYS:HB2	3:A:68:LYS:HZ2	1.74	0.52
3:A:174:ILE:O	3:A:195:LEU:HD12	2.10	0.52
3:A:46:ILE:HB	3:A:53:ILE:CD1	2.40	0.52
3:A:11:LEU:HD23	3:A:11:LEU:N	2.20	0.52
3:A:119:ILE:HG23	3:A:124:ASP:CB	2.40	0.52
3:A:164:ASN:O	3:A:168:LYS:HG2	2.09	0.51
3:A:177:VAL:HG11	3:A:191:MET:HE2	1.91	0.51
3:A:18:MET:HG3	3:A:22:LEU:CD2	2.40	0.51
3:A:26:GLU:O	3:A:30:SER:O	2.29	0.51
3:A:152:ARG:HA	3:A:155:MET:HB2	1.92	0.51
3:A:182:ARG:HH12	3:A:273:THR:HG21	1.74	0.51
3:A:46:ILE:HG13	3:A:47:ALA:N	2.22	0.51
3:A:255:ILE:HG12	3:A:256:ASP:N	2.26	0.51
3:A:124:ASP:O	3:A:128:ASN:ND2	2.43	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:127:LYS:HB2	3:A:128:ASN:HD21	1.74	0.50
3:A:31:GLN:HB2	3:A:112:ARG:NH1	2.24	0.50
3:A:230:LYS:HG3	3:A:235:PHE:HD2	1.75	0.50
3:A:298:ILE:O	3:A:311:LEU:HB2	2.11	0.50
3:A:248:LYS:O	3:A:248:LYS:HG2	2.11	0.50
3:A:73:ILE:CG2	3:A:77:LEU:HD21	2.40	0.50
3:A:244:LYS:C	3:A:245:ASN:HD22	2.17	0.50
3:A:272:PHE:CD1	6:A:338:TTP:H4'	2.41	0.50
3:A:286:ALA:HB1	3:A:291:PHE:HB2	1.93	0.50
2:P:6:DT:C2'	2:P:7:DG:H5''	2.41	0.50
3:A:18:MET:CE	3:A:82:LEU:HD13	2.42	0.50
3:A:165:GLU:HB3	3:A:217:GLN:HG3	1.93	0.50
2:P:5:DA:H2''	2:P:6:DT:H5'	1.93	0.49
3:A:29:VAL:HG21	3:A:94:SER:CB	2.34	0.49
3:A:244:LYS:HB3	3:A:245:ASN:HD22	1.77	0.49
3:A:278:PHE:HB2	3:A:333:ARG:O	2.12	0.49
3:A:120:LYS:N	3:A:124:ASP:OD2	2.43	0.49
3:A:183:ARG:NH1	3:A:275:SER:HA	2.26	0.49
3:A:270:LEU:HD12	3:A:270:LEU:O	2.12	0.49
3:A:150:ILE:N	3:A:188:SER:O	2.37	0.48
2:P:6:DT:H2''	2:P:7:DG:C8	2.48	0.48
3:A:15:ILE:O	3:A:19:LEU:HB2	2.13	0.48
3:A:138:ILE:HD11	3:A:226:ASP:HB3	1.94	0.48
3:A:295:GLU:OE1	3:A:295:GLU:N	2.44	0.48
3:A:42:ALA:O	3:A:46:ILE:HG23	2.13	0.48
3:A:59:ALA:O	3:A:62:LEU:HB2	2.13	0.48
3:A:286:ALA:O	3:A:291:PHE:HB2	2.14	0.48
3:A:295:GLU:CD	3:A:295:GLU:H	2.22	0.48
3:A:243:SER:OG	3:A:249:GLU:HA	2.14	0.48
3:A:327:TYR:CD1	3:A:328:ARG:N	2.80	0.48
3:A:81:LYS:NZ	3:A:86:GLU:HB3	2.29	0.47
3:A:264:GLN:NE2	3:A:296:TYR:HB3	2.29	0.47
3:A:298:ILE:N	7:A:578:HOH:O	2.33	0.47
3:A:155:MET:HE3	3:A:181:PHE:HB2	1.96	0.47
2:P:5:DA:O5'	3:A:107:GLY:HA3	2.15	0.47
3:A:18:MET:HE2	3:A:82:LEU:CD2	2.39	0.47
3:A:306:VAL:HG23	3:A:307:ALA:N	2.30	0.47
3:A:293:ILE:CD1	3:A:298:ILE:HG13	2.21	0.47
3:A:328:ARG:O	3:A:333:ARG:NE	2.39	0.47
2:P:4:DA:H5'	7:P:569:HOH:O	2.14	0.47
3:A:155:MET:SD	3:A:158:MET:HE3	2.54	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:195:LEU:HD12	3:A:195:LEU:HA	1.55	0.47
3:A:296:TYR:O	3:A:297:THR:HG23	2.15	0.47
3:A:299:ARG:HG2	3:A:310:PRO:N	2.29	0.47
3:A:180:SER:HA	3:A:183:ARG:HB2	1.96	0.47
3:A:272:PHE:O	6:A:338:TTP:O3'	2.32	0.47
3:A:287:LEU:HA	3:A:287:LEU:HD13	1.39	0.47
3:A:270:LEU:HD23	3:A:319:ILE:HD13	1.96	0.46
3:A:100:LEU:HD21	3:A:119:ILE:O	2.15	0.46
3:A:228:LEU:HD12	3:A:228:LEU:HA	1.57	0.46
3:A:99:PHE:CD1	3:A:99:PHE:C	2.94	0.46
3:A:207:GLN:O	3:A:210:LEU:HB2	2.15	0.46
3:A:254:ARG:CZ	3:A:254:ARG:HB3	2.46	0.46
3:A:282:MET:HE3	7:A:555:HOH:O	2.14	0.46
3:A:194:LEU:N	3:A:194:LEU:CD1	2.78	0.46
3:A:270:LEU:CD1	3:A:333:ARG:HH12	2.28	0.46
3:A:18:MET:HE2	3:A:82:LEU:HD13	1.97	0.45
3:A:156:LEU:HA	3:A:156:LEU:HD23	1.38	0.45
3:A:298:ILE:O	3:A:298:ILE:HG23	2.15	0.45
3:A:196:THR:OG1	3:A:262:LYS:HA	2.16	0.45
3:A:204:SER:O	3:A:204:SER:OG	2.28	0.45
3:A:277:ILE:HD13	3:A:277:ILE:N	2.07	0.45
3:A:279:ASN:HD22	3:A:279:ASN:HA	1.51	0.45
3:A:270:LEU:HD21	3:A:282:MET:HE1	1.98	0.45
3:A:15:ILE:HB	3:A:46:ILE:CD1	2.47	0.45
3:A:104:SER:N	7:A:600:HOH:O	2.28	0.45
3:A:212:HIS:N	3:A:212:HIS:CD2	2.84	0.45
3:A:31:GLN:HE21	3:A:112:ARG:NH1	2.14	0.45
3:A:245:ASN:ND2	3:A:245:ASN:H	2.05	0.45
3:A:259:LEU:HD12	3:A:259:LEU:HA	1.58	0.45
3:A:278:PHE:CD2	3:A:333:ARG:HB3	2.52	0.45
3:A:145:ASP:OD2	3:A:252:HIS:HB2	2.17	0.44
3:A:272:PHE:C	6:A:338:TTP:HO3'	2.25	0.44
3:A:56:GLY:O	3:A:59:ALA:HB3	2.16	0.44
3:A:62:LEU:HD12	3:A:62:LEU:HA	1.59	0.44
3:A:253:ARG:HG3	3:A:253:ARG:HH11	1.82	0.44
3:A:44:SER:O	3:A:47:ALA:HB3	2.17	0.44
3:A:163:LEU:HD21	3:A:175:ALA:HB3	1.99	0.44
3:A:286:ALA:CA	3:A:323:ILE:HG21	2.48	0.44
3:A:169:VAL:HG22	3:A:173:TYR:HE2	1.83	0.44
3:A:172:GLU:HG2	3:A:198:PRO:HG2	1.99	0.44
3:A:182:ARG:NH1	3:A:273:THR:CG2	2.80	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:309:GLU:HA	3:A:310:PRO:HD2	1.53	0.44
3:A:58:GLU:O	3:A:61:LYS:HG3	2.18	0.43
3:A:59:ALA:C	3:A:65:VAL:HG21	2.44	0.43
3:A:100:LEU:HD12	3:A:100:LEU:HA	1.62	0.43
3:A:146:PHE:HD1	3:A:146:PHE:HA	1.59	0.43
3:A:157:GLN:O	3:A:160:ASP:HB3	2.18	0.43
3:A:205:THR:HA	7:A:624:HOH:O	2.19	0.43
2:P:6:DT:H6	2:P:6:DT:H2'	1.16	0.43
3:A:105:GLY:HA3	3:A:136:GLN:HG2	2.00	0.43
3:A:165:GLU:H	3:A:165:GLU:HG2	1.50	0.43
3:A:266:TYR:HA	3:A:269:VAL:HB	1.98	0.43
3:A:228:LEU:HB2	3:A:236:MET:O	2.19	0.43
3:A:327:TYR:CD1	3:A:327:TYR:C	2.96	0.43
2:P:1:DT:H2'	2:P:2:DC:C6	2.53	0.43
3:A:11:LEU:CD2	3:A:11:LEU:N	2.79	0.43
2:P:2:DC:H2'	2:P:2:DC:O5'	2.19	0.42
3:A:99:PHE:HD1	3:A:100:LEU:HD13	1.84	0.42
3:A:217:GLN:NE2	3:A:217:GLN:HA	2.33	0.42
1:T:7:DA:H1'	7:T:634:HOH:O	2.19	0.42
3:A:79:THR:C	3:A:81:LYS:H	2.24	0.42
3:A:163:LEU:HA	3:A:163:LEU:HD23	1.07	0.42
3:A:277:ILE:HG12	3:A:335:GLU:CA	2.43	0.42
3:A:277:ILE:CG1	3:A:335:GLU:CB	2.98	0.42
3:A:286:ALA:HB2	3:A:293:ILE:HD11	1.99	0.42
3:A:291:PHE:CD1	3:A:300:PRO:HA	2.54	0.42
3:A:26:GLU:HA	3:A:30:SER:OG	2.20	0.42
3:A:303:VAL:C	3:A:305:GLY:N	2.77	0.42
2:P:5:DA:P	3:A:107:GLY:HA3	2.60	0.42
3:A:249:GLU:HG3	3:A:250:TYR:N	2.35	0.42
3:A:288:GLU:C	3:A:290:GLY:H	2.26	0.42
6:A:338:TTP:H3'	7:A:631:HOH:O	2.19	0.42
2:P:5:DA:H2'	2:P:6:DT:H71	2.02	0.42
3:A:235:PHE:CD1	3:A:235:PHE:C	2.97	0.42
3:A:22:LEU:HD13	3:A:85:LEU:CD1	2.50	0.42
3:A:54:LYS:O	3:A:55:SER:HB3	2.20	0.42
3:A:299:ARG:NH2	7:A:580:HOH:O	2.44	0.42
3:A:15:ILE:HD11	3:A:77:LEU:HD11	2.02	0.41
3:A:218:LEU:N	3:A:218:LEU:CD1	2.83	0.41
3:A:133:ASN:OD1	3:A:136:GLN:HG3	2.19	0.41
3:A:271:TYR:CD2	3:A:272:PHE:N	2.88	0.41
3:A:332:ASP:C	3:A:334:SER:H	2.28	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:59:ALA:O	3:A:65:VAL:HG21	2.21	0.41
3:A:69:ILE:O	3:A:72:LYS:N	2.53	0.41
3:A:207:GLN:HB3	3:A:210:LEU:HG	2.03	0.41
3:A:321:ASP:O	3:A:324:GLN:N	2.52	0.41
3:A:48:LYS:O	3:A:50:PRO:HD3	2.20	0.41
3:A:60:LYS:O	3:A:60:LYS:HG3	2.21	0.41
3:A:316:GLU:O	3:A:320:PHE:CD2	2.74	0.41
3:A:77:LEU:HD13	3:A:77:LEU:N	2.35	0.41
3:A:277:ILE:HG13	3:A:335:GLU:CB	2.51	0.41
3:A:285:HIS:C	3:A:287:LEU:N	2.78	0.41
3:A:299:ARG:CD	3:A:310:PRO:HD3	2.51	0.41
1:T:4:DT:H1'	7:T:527:HOH:O	2.20	0.41
3:A:34:HIS:O	3:A:38:ALA:N	2.53	0.41
3:A:121:THR:OG1	3:A:123:GLU:N	2.52	0.41
3:A:122:LEU:CD2	3:A:126:ARG:CZ	2.99	0.41
3:A:182:ARG:HH11	3:A:182:ARG:CG	2.23	0.41
3:A:251:PRO:O	3:A:253:ARG:HD2	2.21	0.41
3:A:103:VAL:HB	3:A:106:ILE:HD12	2.02	0.41
3:A:155:MET:HE3	3:A:181:PHE:CB	2.51	0.41
3:A:165:GLU:O	3:A:168:LYS:N	2.54	0.41
3:A:30:SER:HB3	7:A:561:HOH:O	2.20	0.40
3:A:127:LYS:HG2	3:A:127:LYS:H	1.44	0.40
3:A:317:LYS:O	3:A:320:PHE:N	2.53	0.40
3:A:27:LYS:HG2	3:A:36:TYR:CD1	2.56	0.40
3:A:322:TYR:N	3:A:322:TYR:CD1	2.84	0.40
3:A:266:TYR:O	3:A:270:LEU:N	2.46	0.40
3:A:16:THR:HG23	3:A:46:ILE:HD11	2.03	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
3	A	325/335 (97%)	266 (82%)	39 (12%)	20 (6%)	1 6

All (20) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	A	185	ALA
3	A	204	SER
3	A	205	THR
3	A	206	LYS
3	A	244	LYS
3	A	246	ASP
3	A	295	GLU
3	A	202	SER
3	A	265	TYR
3	A	91	ASP
3	A	222	HIS
3	A	286	ALA
3	A	310	PRO
3	A	13	GLY
3	A	80	GLY
3	A	207	GLN
3	A	247	GLU
3	A	304	THR
3	A	309	GLU
3	A	316	GLU

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
3	A	288/295 (98%)	212 (74%)	76 (26%)	0 3

All (76) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	A	10	THR

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Mol	Chain	Res	Type
3	A	18	MET
3	A	19	LEU
3	A	21	GLU
3	A	22	LEU
3	A	27	LYS
3	A	30	SER
3	A	33	ILE
3	A	37	ASN
3	A	41	LYS
3	A	46	ILE
3	A	52	LYS
3	A	62	LEU
3	A	65	VAL
3	A	67	THR
3	A	68	LYS
3	A	72	LYS
3	A	77	LEU
3	A	79	THR
3	A	85	LEU
3	A	89	ARG
3	A	92	ASP
3	A	93	THR
3	A	94	SER
3	A	97	ILE
3	A	100	LEU
3	A	101	THR
3	A	115	VAL
3	A	121	THR
3	A	128	ASN
3	A	138	ILE
3	A	152	ARG
3	A	153	GLU
3	A	157	GLN
3	A	161	ILE
3	A	168	LYS
3	A	169	VAL
3	A	171	SER
3	A	174	ILE
3	A	182	ARG
3	A	194	LEU
3	A	205	THR
3	A	213	GLN

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Mol	Chain	Res	Type
3	A	214	VAL
3	A	218	LEU
3	A	225	THR
3	A	226	ASP
3	A	228	LEU
3	A	229	SER
3	A	230	LYS
3	A	245	ASN
3	A	248	LYS
3	A	249	GLU
3	A	253	ARG
3	A	255	ILE
3	A	256	ASP
3	A	258	ARG
3	A	264	GLN
3	A	270	LEU
3	A	277	ILE
3	A	279	ASN
3	A	287	LEU
3	A	289	LYS
3	A	293	ILE
3	A	294	ASN
3	A	295	GLU
3	A	298	ILE
3	A	301	LEU
3	A	306	VAL
3	A	309	GLU
3	A	311	LEU
3	A	314	ASP
3	A	324	GLN
3	A	325	TRP
3	A	331	LYS
3	A	335	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
3	A	12	ASN
3	A	28	ASN
3	A	31	GLN
3	A	128	ASN
3	A	159	GLN

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Mol	Chain	Res	Type
3	A	212	HIS
3	A	213	GLN
3	A	217	GLN
3	A	245	ASN
3	A	279	ASN
3	A	324	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 5 ligands modelled in this entry, 4 are monoatomic - leaving 1 for Mogul analysis.

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	$\# Z > 2$
6	TTP	A	338	4	19,20,30	1.31	1 (5%)	25,31,47	1.42	1 (4%)

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
6	TTP	A	338	4	-	10/18/28/34	0/1/1/2

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	338	TTP	PB-O3A	4.27	1.64	1.59

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	338	TTP	O3'-C3'-C2'	-5.24	99.14	111.43

There are no chirality outliers.

All (10) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	A	338	TTP	C5'-O5'-PA-O1A
6	A	338	TTP	C5'-O5'-PA-O2A
6	A	338	TTP	C5'-O5'-PA-O3A
6	A	338	TTP	PB-O3B-PG-O2G
6	A	338	TTP	O4'-C4'-C5'-O5'
6	A	338	TTP	PA-O3A-PB-O2B
6	A	338	TTP	PG-O3B-PB-O1B
6	A	338	TTP	PG-O3B-PB-O2B
6	A	338	TTP	PB-O3B-PG-O1G
6	A	338	TTP	PA-O3A-PB-O1B

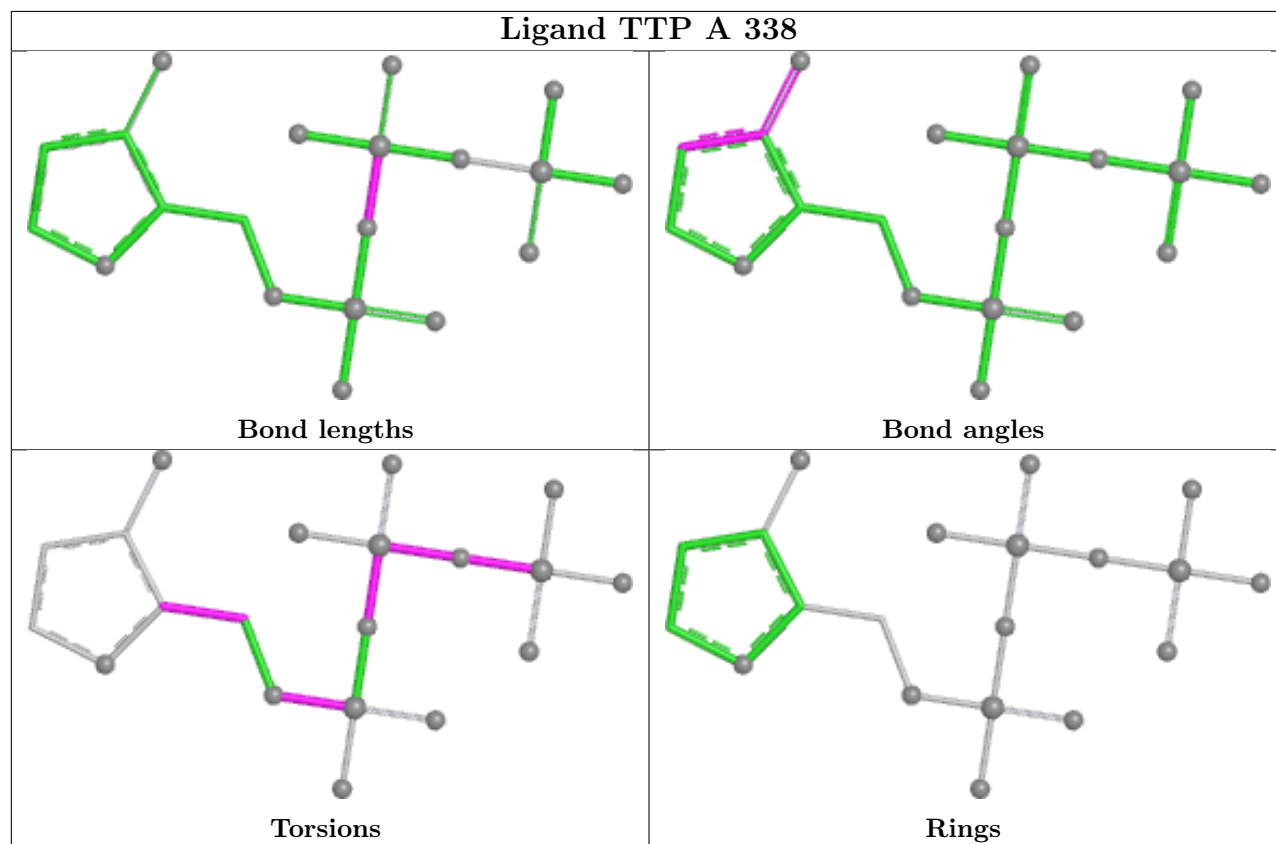
There are no ring outliers.

1 monomer is involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	338	TTP	5	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	T	8/8 (100%)	-0.49	0	100 100	13, 37, 99, 99	0
2	P	7/7 (100%)	-1.02	0	100 100	13, 25, 37, 64	0
3	A	325/335 (97%)	-0.94	1 (0%)	90 79	2, 32, 86, 100	0
All	All	340/350 (97%)	-0.93	1 (0%)	90 79	2, 32, 88, 100	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	A	246	ASP	2.4

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 oligosaccharides 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.

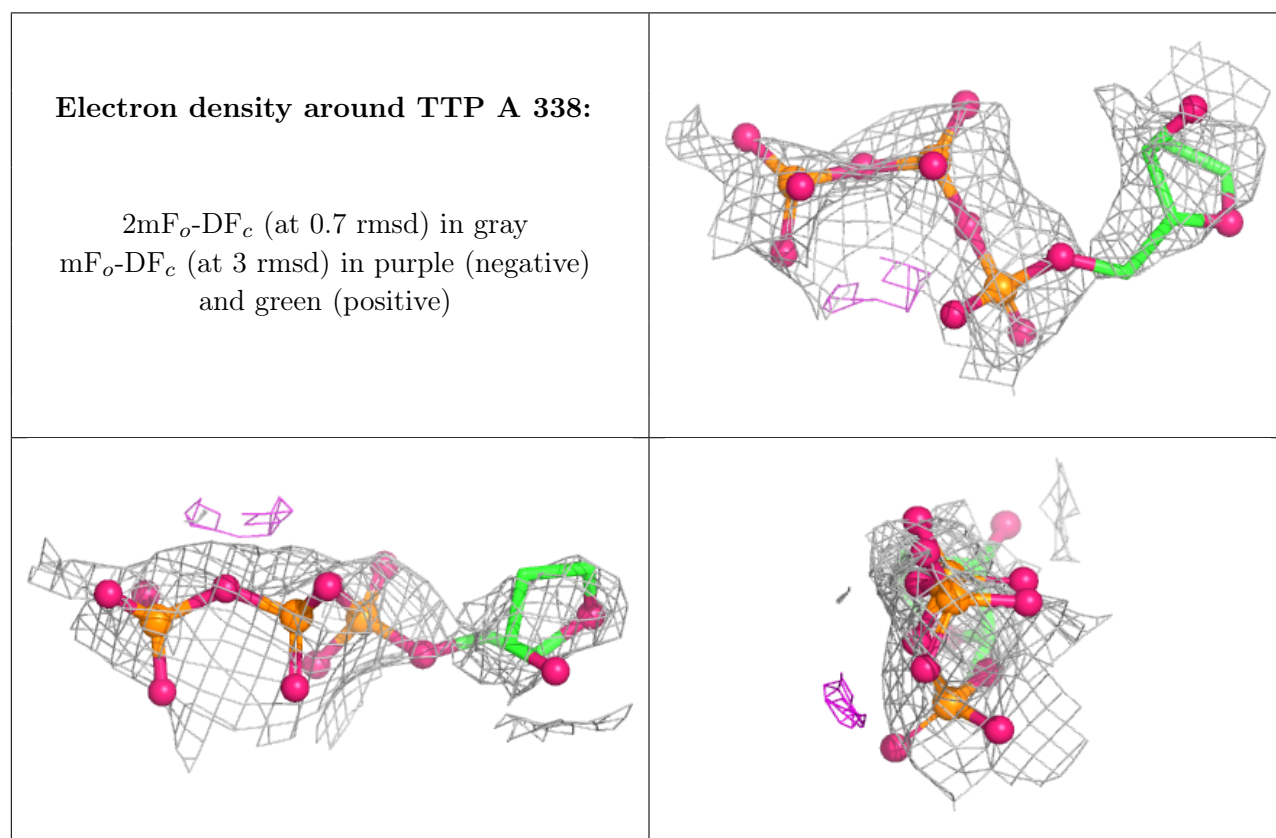
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	MN	A	340	1/1	0.88	0.06	30,30,30,30	1
4	MN	A	339	1/1	0.89	0.05	30,30,30,30	1

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
6	TTP	A	338	20/29	0.89	0.11	69,72,78,80	20
5	NA	A	342	1/1	0.95	0.10	44,44,44,44	0
5	NA	A	341	1/1	0.99	0.03	6,6,6,6	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.



6.5 Other polymers [i](#)

There are no such residues in this entry.