



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 17, 2026 – 10:09 PM UTC

PDB ID : 1ZM7 / pdb_00001zm7
Title : Crystal structure of D. melanogaster deoxyribonucleoside kinase mutant N64D in complex with dTTP
Authors : Welin, M.; Skovgaard, T.; Knecht, W.; Berenstein, D.; Munch-Petersen, B.; Piskur, J.; Eklund, H.
Deposited on : 2005-05-10
Resolution : 2.20 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

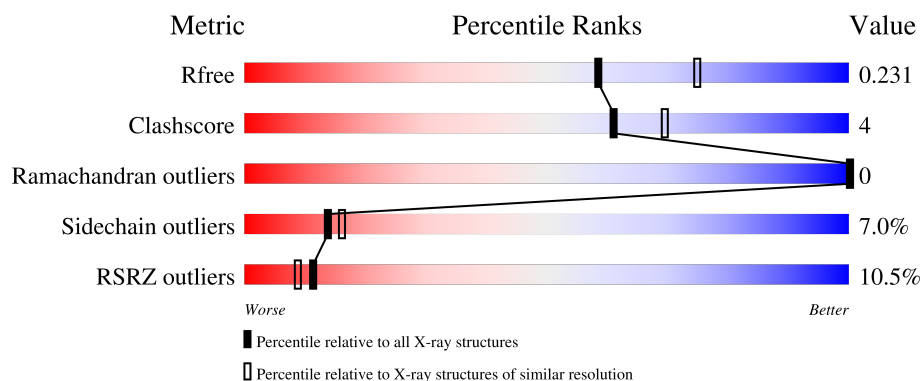
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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}	180053	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	A	230	8% 75% 9% 14%
1	B	230	8% 75% 10% 14%
1	C	230	9% 74% 10% 14%
1	D	230	10% 77% 7% 14%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6887 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.

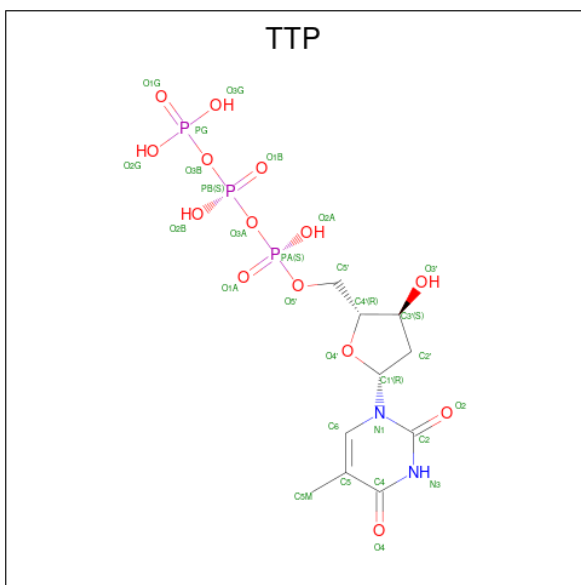
- Molecule 1 is a protein called Deoxynucleoside kinase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	197	Total	C	N	O	S	0	0	0
			1640	1046	281	303	10			
1	B	198	Total	C	N	O	S	0	0	0
			1648	1052	282	304	10			
1	C	198	Total	C	N	O	S	0	0	0
			1648	1052	282	304	10			
1	D	197	Total	C	N	O	S	0	0	0
			1640	1046	281	303	10			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	64	ASP	ASN	engineered mutation	UNP Q9XZT6
B	64	ASP	ASN	engineered mutation	UNP Q9XZT6
C	64	ASP	ASN	engineered mutation	UNP Q9XZT6
D	64	ASP	ASN	engineered mutation	UNP Q9XZT6

- Molecule 2 is THYMIDINE-5'-TRIPHOSPHATE (CCD ID: TTP) (formula: $C_{10}H_{17}N_2O_{14}P_3$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			29	10	2	14	3		
2	B	1	Total	C	N	O	P	0	0
			29	10	2	14	3		
2	C	1	Total	C	N	O	P	0	0
			29	10	2	14	3		
2	D	1	Total	C	N	O	P	0	0
			29	10	2	14	3		

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	D	1	Total	Mg	0	0
			1	1		

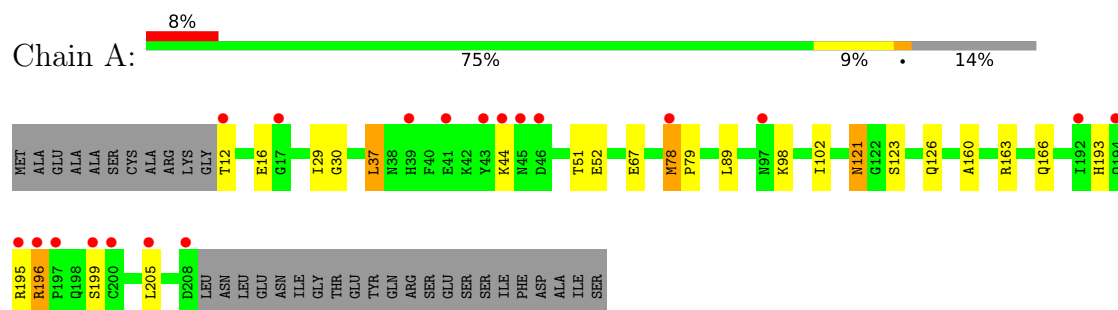
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	38	Total	O	0	0
			38	38		
4	B	50	Total	O	0	0
			50	50		
4	C	53	Total	O	0	0
			53	53		
4	D	53	Total	O	0	0
			53	53		

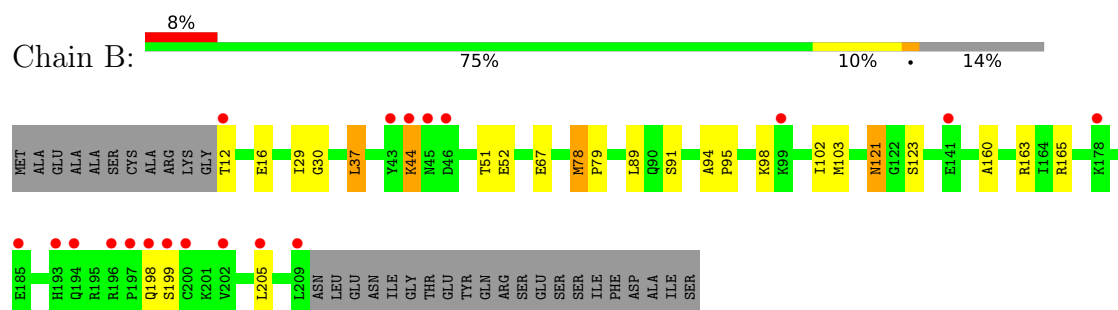
3 Residue-property plots [i](#)

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.

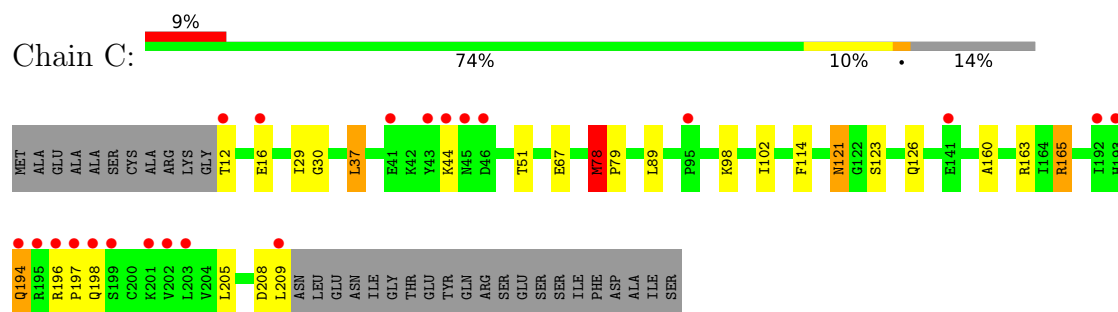
• Molecule 1: Deoxynucleoside kinase



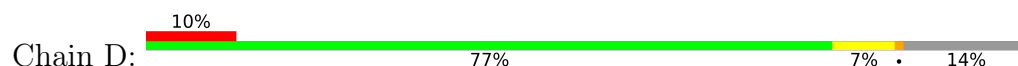
• Molecule 1: Deoxynucleoside kinase

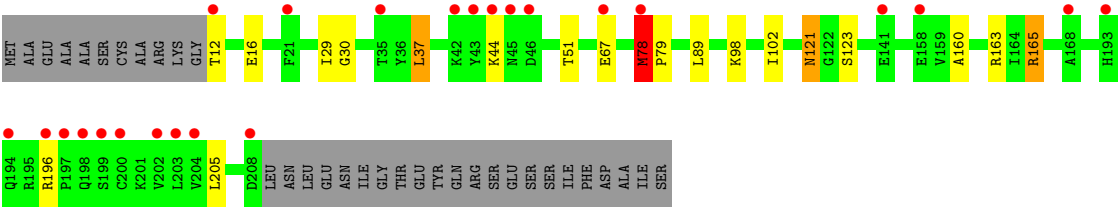


• Molecule 1: Deoxynucleoside kinase



• Molecule 1: Deoxynucleoside kinase





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	67.04Å 119.27Å 68.39Å 90.00° 92.59° 90.00°	Depositor
Resolution (Å)	44.72 – 2.20 44.72 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.9 (44.72-2.20) 99.9 (44.72-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.95 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.1.24	Depositor
R, R_{free}	0.212 , 0.237 0.211 , 0.231	Depositor DCC
R_{free} test set	2763 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	28.5	Xtriage
Anisotropy	0.762	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 37.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.000 for l,k,-h 0.036 for h,-k,-l 0.000 for l,-k,h	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6887	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 64.40 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 8.3320e-06. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: MG, TTP

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	0.66	0/1678	0.80	0/2268
1	B	0.67	1/1686 (0.1%)	0.80	0/2279
1	C	0.66	1/1686 (0.1%)	0.78	0/2279
1	D	0.66	1/1678 (0.1%)	0.79	0/2268
All	All	0.66	3/6728 (0.0%)	0.79	0/9094

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	52	GLU	CA-C	5.99	1.57	1.53
1	C	78	MET	N-CA	5.40	1.50	1.46
1	D	78	MET	N-CA	5.07	1.50	1.46

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

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	1640	0	1627	17	0
1	B	1648	0	1638	18	0
1	C	1648	0	1638	22	0
1	D	1640	0	1626	13	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	29	0	13	0	0
2	B	29	0	13	0	0
2	C	29	0	13	1	0
2	D	29	0	13	0	0
3	D	1	0	0	0	0
4	A	38	0	0	0	0
4	B	50	0	0	1	0
4	C	53	0	0	0	0
4	D	53	0	0	0	0
All	All	6887	0	6581	56	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 4.

All (56) 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:GLN:HE22	1:C:126:GLN:HE22	1.07	0.94
1:B:91:SER:HB3	1:B:103:MET:HE3	1.60	0.84
1:A:126:GLN:NE2	1:C:126:GLN:HE22	1.75	0.83
1:A:126:GLN:HE22	1:C:126:GLN:NE2	1.75	0.83
1:A:78:MET:SD	1:B:79:PRO:HB3	2.21	0.81
1:C:78:MET:SD	1:D:79:PRO:HB3	2.26	0.76
1:C:79:PRO:HB3	1:D:78:MET:SD	2.27	0.74
1:B:91:SER:CB	1:B:103:MET:HE3	2.24	0.66
1:A:79:PRO:HB3	1:B:78:MET:SD	2.38	0.64
1:B:121:ASN:C	1:B:121:ASN:HD22	2.10	0.60
1:D:121:ASN:C	1:D:121:ASN:HD22	2.11	0.58
1:A:121:ASN:HD22	1:A:121:ASN:C	2.11	0.58
1:C:121:ASN:C	1:C:121:ASN:HD22	2.11	0.57
1:C:165:ARG:HH11	1:C:165:ARG:CG	2.18	0.57
1:C:37:LEU:HD22	1:C:51:THR:HG22	1.88	0.56
1:D:37:LEU:HD22	1:D:51:THR:HG22	1.88	0.56
1:B:37:LEU:HD22	1:B:51:THR:HG22	1.88	0.55
1:A:78:MET:SD	1:B:79:PRO:CB	2.93	0.55
1:A:37:LEU:HD22	1:A:51:THR:HG22	1.88	0.54
1:C:165:ARG:HH11	1:C:165:ARG:HG3	1.71	0.54
1:B:198:GLN:HA	1:B:198:GLN:OE1	2.06	0.54
1:A:193:HIS:HB3	1:A:195:ARG:CZ	2.38	0.54
1:B:44:LYS:HD3	4:B:341:HOH:O	2.08	0.53
1:C:165:ARG:HG3	1:C:165:ARG:NH1	2.23	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:79:PRO:CB	1:D:78:MET:SD	2.99	0.50
1:C:78:MET:SD	1:D:79:PRO:CB	3.00	0.48
1:D:30:GLY:O	1:D:163:ARG:HD2	2.15	0.47
1:B:30:GLY:O	1:B:163:ARG:HD2	2.15	0.47
1:A:30:GLY:O	1:A:163:ARG:HD2	2.14	0.47
1:C:30:GLY:O	1:C:163:ARG:HD2	2.15	0.46
1:A:121:ASN:HD22	1:A:123:SER:H	1.64	0.46
1:C:121:ASN:HD22	1:C:123:SER:H	1.64	0.46
1:A:78:MET:SD	1:B:79:PRO:CA	3.05	0.45
1:D:121:ASN:HD22	1:D:123:SER:H	1.64	0.44
1:B:121:ASN:HD22	1:B:123:SER:H	1.64	0.44
1:B:121:ASN:ND2	1:B:123:SER:H	2.16	0.44
1:A:196:ARG:HD3	1:A:196:ARG:HA	1.78	0.44
1:D:121:ASN:ND2	1:D:123:SER:H	2.16	0.44
1:B:44:LYS:HE3	1:C:194:GLN:HG2	2.00	0.43
1:C:121:ASN:ND2	1:C:123:SER:H	2.17	0.43
1:A:121:ASN:ND2	1:A:123:SER:H	2.17	0.42
1:C:29:ILE:O	1:C:160:ALA:HB1	2.20	0.42
1:D:37:LEU:HD21	1:D:102:ILE:HG22	2.02	0.42
1:A:29:ILE:O	1:A:160:ALA:HB1	2.20	0.41
1:C:79:PRO:CA	1:D:78:MET:SD	3.08	0.41
1:B:37:LEU:HD21	1:B:102:ILE:HG22	2.02	0.41
1:D:29:ILE:O	1:D:160:ALA:HB1	2.20	0.41
1:C:37:LEU:HD21	1:C:102:ILE:HG22	2.03	0.41
1:C:196:ARG:HA	1:C:197:PRO:HD3	1.91	0.41
1:C:208:ASP:O	1:C:209:LEU:HB2	2.20	0.41
1:D:165:ARG:HD2	1:D:165:ARG:HA	1.76	0.41
1:A:37:LEU:HD21	1:A:102:ILE:HG22	2.03	0.41
1:B:94:ALA:HA	1:B:95:PRO:HD3	1.88	0.41
1:A:78:MET:SD	1:B:79:PRO:HA	2.60	0.41
1:B:29:ILE:O	1:B:160:ALA:HB1	2.21	0.40
1:C:114:PHE:CG	2:C:302:TTP:C4	3.10	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	
1	A	195/230 (85%)	188 (96%)	7 (4%)	0	100	100
1	B	196/230 (85%)	189 (96%)	7 (4%)	0	100	100
1	C	196/230 (85%)	189 (96%)	7 (4%)	0	100	100
1	D	195/230 (85%)	188 (96%)	7 (4%)	0	100	100
All	All	782/920 (85%)	754 (96%)	28 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	182/208 (88%)	168 (92%)	14 (8%)	12	13
1	B	183/208 (88%)	171 (93%)	12 (7%)	15	18
1	C	183/208 (88%)	170 (93%)	13 (7%)	13	16
1	D	182/208 (88%)	170 (93%)	12 (7%)	15	18
All	All	730/832 (88%)	679 (93%)	51 (7%)	14	16

All (51) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	12	THR
1	A	16	GLU
1	A	37	LEU
1	A	44	LYS
1	A	52	GLU
1	A	67	GLU
1	A	78	MET
1	A	89	LEU

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Mol	Chain	Res	Type
1	A	98	LYS
1	A	121	ASN
1	A	166	GLN
1	A	196	ARG
1	A	199	SER
1	A	205	LEU
1	B	12	THR
1	B	16	GLU
1	B	37	LEU
1	B	44	LYS
1	B	67	GLU
1	B	78	MET
1	B	89	LEU
1	B	98	LYS
1	B	121	ASN
1	B	165	ARG
1	B	199	SER
1	B	205	LEU
1	C	12	THR
1	C	16	GLU
1	C	37	LEU
1	C	44	LYS
1	C	67	GLU
1	C	78	MET
1	C	89	LEU
1	C	98	LYS
1	C	121	ASN
1	C	165	ARG
1	C	194	GLN
1	C	198	GLN
1	C	205	LEU
1	D	12	THR
1	D	16	GLU
1	D	37	LEU
1	D	44	LYS
1	D	67	GLU
1	D	78	MET
1	D	89	LEU
1	D	98	LYS
1	D	121	ASN
1	D	165	ARG
1	D	196	ARG

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Mol	Chain	Res	Type
1	D	205	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	121	ASN
1	A	130	ASN
1	B	45	ASN
1	B	121	ASN
1	B	126	GLN
1	B	184	HIS
1	C	121	ASN
1	C	126	GLN
1	D	45	ASN
1	D	121	ASN
1	D	126	GLN
1	D	184	HIS

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, 1 is monoatomic - leaving 4 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	$\# Z > 2$
2	TTP	B	301	-	29,30,30	2.04	8 (27%)	43,47,47	1.88	10 (23%)
2	TTP	A	300	-	29,30,30	2.04	9 (31%)	43,47,47	1.78	8 (18%)
2	TTP	C	302	-	29,30,30	2.09	10 (34%)	43,47,47	2.05	8 (18%)
2	TTP	D	303	3	29,30,30	2.11	7 (24%)	43,47,47	1.83	10 (23%)

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
2	TTP	B	301	-	-	3/22/34/34	0/2/2/2
2	TTP	A	300	-	-	2/22/34/34	0/2/2/2
2	TTP	C	302	-	-	2/22/34/34	0/2/2/2
2	TTP	D	303	3	-	2/22/34/34	0/2/2/2

All (34) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	301	TTP	C1'-N1	-4.96	1.35	1.48
2	D	303	TTP	C3'-C4'	-4.86	1.40	1.53
2	C	302	TTP	C1'-N1	-4.71	1.36	1.48
2	A	300	TTP	C3'-C4'	-4.58	1.40	1.53
2	C	302	TTP	C2'-C3'	-4.03	1.42	1.52
2	C	302	TTP	C3'-C4'	-4.03	1.42	1.53
2	D	303	TTP	C2'-C3'	-4.02	1.42	1.52
2	B	301	TTP	C2'-C3'	-3.83	1.43	1.52
2	C	302	TTP	C2'-C1'	-3.80	1.42	1.52
2	B	301	TTP	C2'-C1'	-3.79	1.42	1.52
2	D	303	TTP	C6-C5	3.76	1.40	1.34
2	A	300	TTP	C2'-C1'	-3.76	1.42	1.52
2	B	301	TTP	C3'-C4'	-3.71	1.43	1.53
2	C	302	TTP	C5'-C4'	-3.70	1.40	1.51
2	D	303	TTP	C4-C5	-3.66	1.38	1.44
2	A	300	TTP	C2'-C3'	-3.64	1.43	1.52
2	A	300	TTP	C1'-N1	-3.62	1.39	1.48
2	D	303	TTP	C2'-C1'	-3.51	1.42	1.52
2	B	301	TTP	C6-C5	3.50	1.40	1.34

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	301	TTP	C5'-C4'	-3.41	1.41	1.51
2	A	300	TTP	C6-C5	3.36	1.40	1.34
2	D	303	TTP	C1'-N1	-3.31	1.39	1.48
2	A	300	TTP	C5'-C4'	-3.26	1.41	1.51
2	D	303	TTP	C5'-C4'	-3.05	1.42	1.51
2	A	300	TTP	PB-O3A	2.84	1.62	1.59
2	C	302	TTP	PA-O3A	2.62	1.62	1.59
2	B	301	TTP	PA-O3A	2.44	1.62	1.59
2	C	302	TTP	C6-C5	2.44	1.38	1.34
2	C	302	TTP	C4-C5	-2.42	1.40	1.44
2	A	300	TTP	C4-C5	-2.26	1.41	1.44
2	C	302	TTP	PB-O3A	2.24	1.61	1.59
2	B	301	TTP	C4-C5	-2.16	1.41	1.44
2	A	300	TTP	PG-O3G	-2.07	1.47	1.54
2	C	302	TTP	PG-O2G	-2.02	1.47	1.54

All (36) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	302	TTP	C4-N3-C2	-5.89	119.62	127.34
2	C	302	TTP	C5-C4-N3	5.85	120.41	115.32
2	B	301	TTP	C5-C4-N3	5.79	120.36	115.32
2	B	301	TTP	C4-N3-C2	-5.02	120.75	127.34
2	C	302	TTP	O4-C4-C5	-4.86	119.36	124.92
2	A	300	TTP	C4-N3-C2	-4.76	121.10	127.34
2	C	302	TTP	N3-C2-N1	4.65	120.94	114.89
2	D	303	TTP	C4-N3-C2	-4.51	121.42	127.34
2	D	303	TTP	C5-C4-N3	4.45	119.19	115.32
2	A	300	TTP	C5-C4-N3	4.33	119.08	115.32
2	A	300	TTP	N3-C2-N1	4.24	120.41	114.89
2	B	301	TTP	N3-C2-N1	4.09	120.21	114.89
2	D	303	TTP	N3-C2-N1	4.07	120.19	114.89
2	D	303	TTP	O4-C4-C5	-3.71	120.67	124.92
2	A	300	TTP	O4-C4-C5	-3.61	120.79	124.92
2	A	300	TTP	O3A-PB-O1B	-3.18	101.13	110.70
2	D	303	TTP	C6-N1-C2	-3.03	118.29	121.30
2	B	301	TTP	C2'-C3'-C4'	2.93	108.74	102.80
2	A	300	TTP	C6-N1-C2	-2.93	118.39	121.30
2	B	301	TTP	O2-C2-N3	-2.83	116.26	121.49
2	C	302	TTP	C6-N1-C2	-2.63	118.68	121.30
2	C	302	TTP	O3A-PB-O1B	-2.61	102.84	110.70
2	D	303	TTP	O3A-PA-O1A	-2.57	102.96	110.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	303	TTP	C2'-C1'-N1	2.53	120.14	113.81
2	D	303	TTP	O2-C2-N3	-2.49	116.90	121.49
2	C	302	TTP	C1'-N1-C2	2.44	122.44	117.66
2	D	303	TTP	C2'-C3'-C4'	2.41	107.69	102.80
2	B	301	TTP	O4-C4-C5	-2.39	122.18	124.92
2	B	301	TTP	O3A-PB-O1B	-2.31	103.77	110.70
2	A	300	TTP	C5M-C5-C4	2.23	121.17	118.78
2	B	301	TTP	O4-C4-N3	-2.18	116.02	120.11
2	C	302	TTP	C2'-C3'-C4'	2.12	107.11	102.80
2	A	300	TTP	O5'-C5'-C4'	2.10	116.15	108.99
2	B	301	TTP	O3A-PA-O1A	-2.09	104.41	110.70
2	B	301	TTP	C5-C6-N1	-2.07	121.07	123.31
2	D	303	TTP	O3A-PB-O1B	-2.02	104.64	110.70

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	301	TTP	C5'-O5'-PA-O1A
2	A	300	TTP	C3'-C4'-C5'-O5'
2	A	300	TTP	O4'-C4'-C5'-O5'
2	C	302	TTP	O4'-C4'-C5'-O5'
2	D	303	TTP	O4'-C4'-C5'-O5'
2	C	302	TTP	C3'-C4'-C5'-O5'
2	B	301	TTP	C3'-C4'-C5'-O5'
2	D	303	TTP	C3'-C4'-C5'-O5'
2	B	301	TTP	O4'-C4'-C5'-O5'

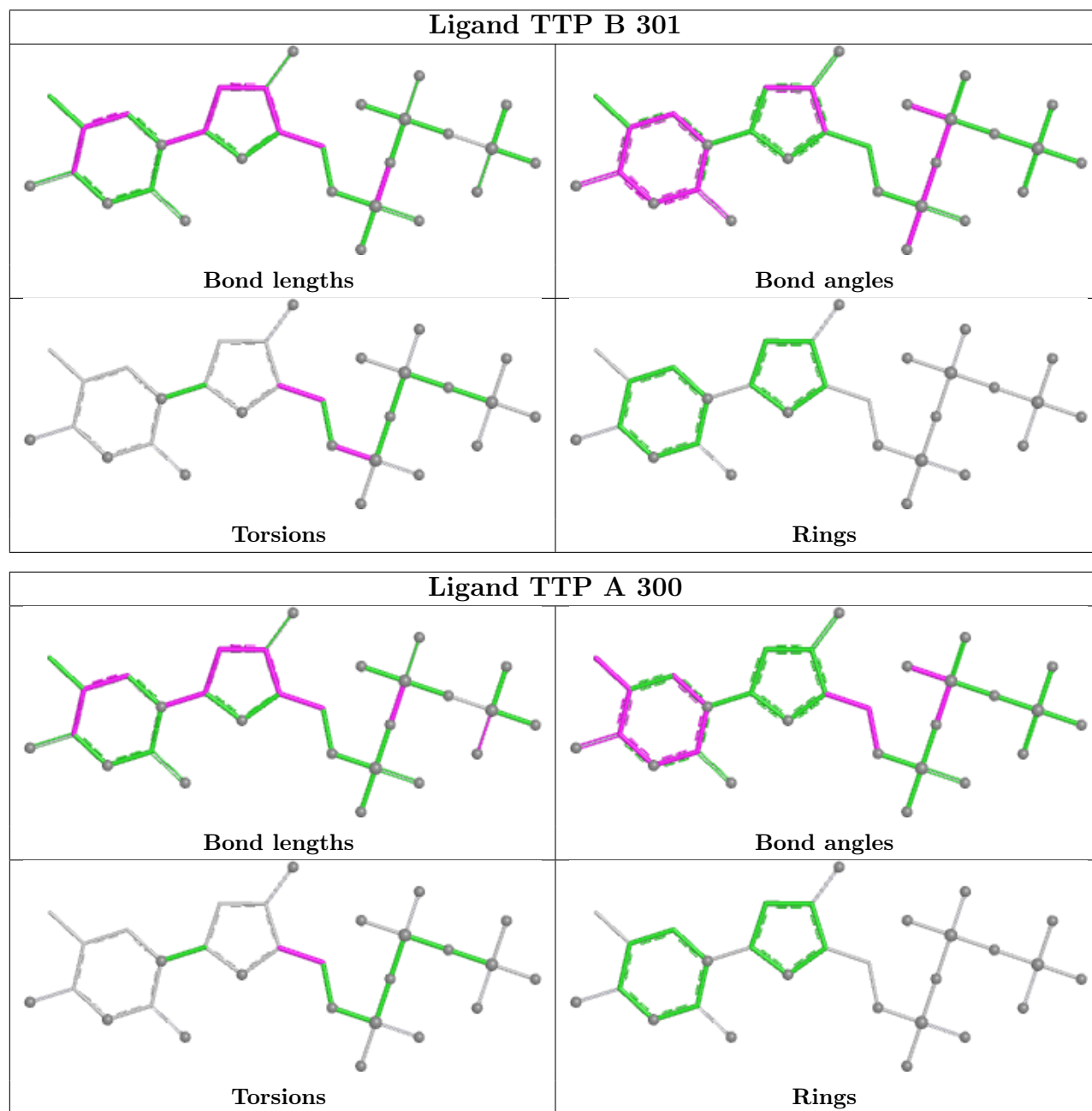
There are no ring outliers.

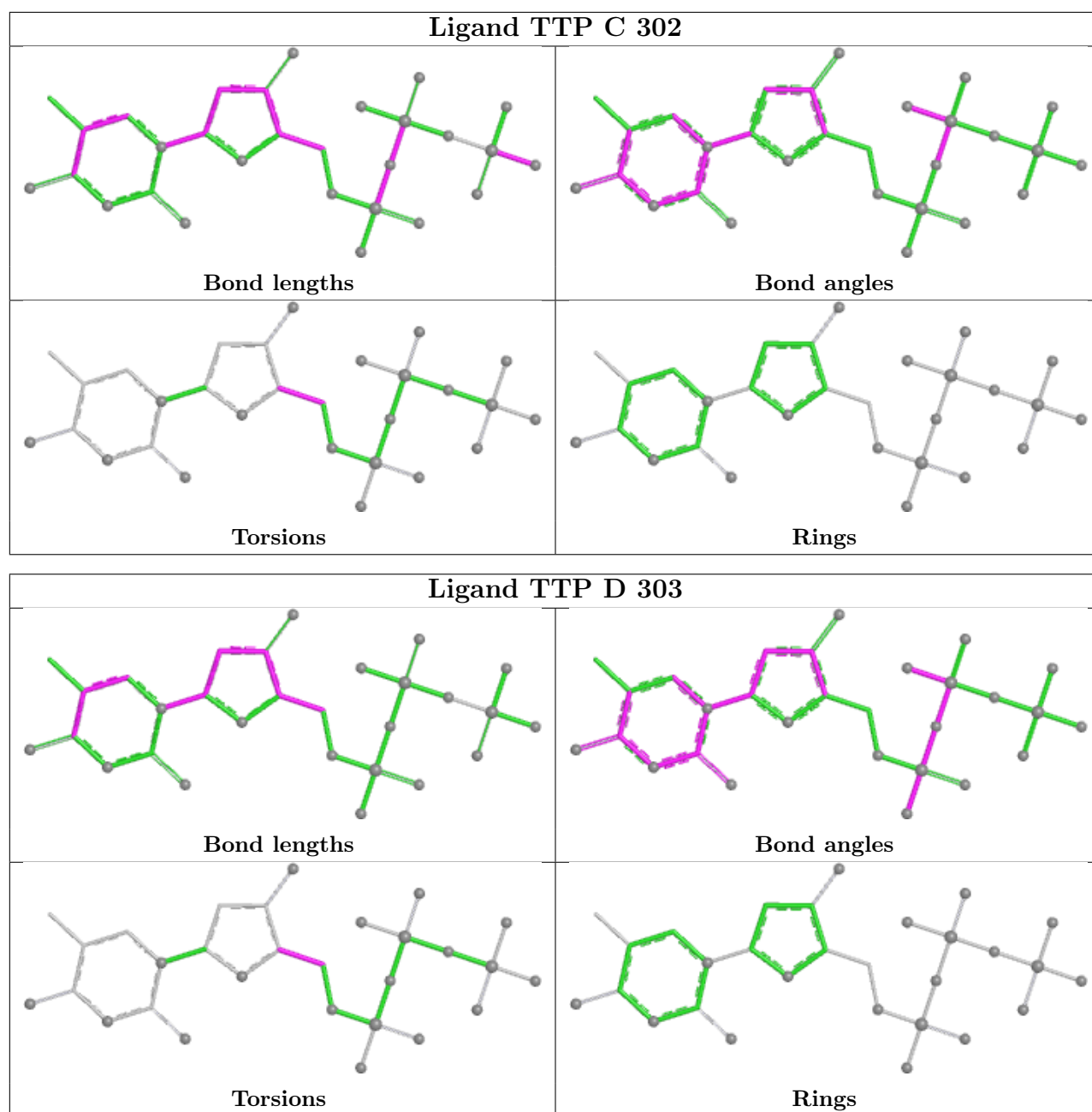
1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	302	TTP	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 [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	A	197/230 (85%)	0.61	19 (9%) 13 11	18, 34, 62, 72	0
1	B	198/230 (86%)	0.58	19 (9%) 13 11	18, 34, 62, 72	0
1	C	198/230 (86%)	0.60	21 (10%) 11 8	18, 34, 62, 72	0
1	D	197/230 (85%)	0.66	24 (12%) 8 6	18, 34, 62, 73	0
All	All	790/920 (85%)	0.61	83 (10%) 11 9	18, 34, 62, 73	0

All (83) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	209	LEU	5.6
1	B	197	PRO	5.3
1	D	197	PRO	5.1
1	D	199	SER	5.1
1	A	199	SER	4.5
1	D	194	GLN	4.4
1	D	43	TYR	4.3
1	C	12	THR	4.2
1	C	196	ARG	4.1
1	A	197	PRO	4.1
1	B	209	LEU	4.0
1	A	196	ARG	3.9
1	C	41	GLU	3.9
1	C	43	TYR	3.8
1	D	196	ARG	3.7
1	A	17	GLY	3.6
1	D	208	ASP	3.4
1	A	45	ASN	3.3
1	A	43	TYR	3.2
1	D	45	ASN	3.2
1	C	194	GLN	3.2

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Mol	Chain	Res	Type	RSRZ
1	C	197	PRO	3.2
1	D	21	PHE	3.1
1	C	46	ASP	3.1
1	B	194	GLN	3.1
1	B	196	ARG	3.1
1	A	195	ARG	3.0
1	B	12	THR	3.0
1	B	199	SER	3.0
1	D	200	CYS	3.0
1	D	198	GLN	3.0
1	C	202	VAL	3.0
1	A	194	GLN	3.0
1	A	208	ASP	2.9
1	D	12	THR	2.9
1	A	44	LYS	2.8
1	C	199	SER	2.8
1	D	141	GLU	2.8
1	B	99	LYS	2.8
1	B	202	VAL	2.8
1	C	44	LYS	2.7
1	A	12	THR	2.7
1	B	198	GLN	2.7
1	B	46	ASP	2.7
1	C	198	GLN	2.6
1	A	41	GLU	2.6
1	C	195	ARG	2.6
1	C	95	PRO	2.5
1	C	16	GLU	2.5
1	C	141	GLU	2.4
1	D	202	VAL	2.4
1	B	185	GLU	2.4
1	D	204	VAL	2.4
1	C	201	LYS	2.4
1	C	192	ILE	2.4
1	D	44	LYS	2.4
1	B	45	ASN	2.3
1	D	67	GLU	2.3
1	C	203	LEU	2.3
1	A	78	MET	2.3
1	B	141	GLU	2.3
1	B	193	HIS	2.3
1	A	46	ASP	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	43	TYR	2.2
1	D	193	HIS	2.2
1	B	44	LYS	2.2
1	B	200	CYS	2.2
1	A	205	LEU	2.2
1	D	46	ASP	2.2
1	A	200	CYS	2.2
1	D	158	GLU	2.2
1	A	192	ILE	2.1
1	B	205	LEU	2.1
1	D	203	LEU	2.1
1	B	178	LYS	2.1
1	A	97	ASN	2.1
1	C	193	HIS	2.1
1	D	168	ALA	2.1
1	D	78	MET	2.0
1	D	35	THR	2.0
1	A	39	HIS	2.0
1	C	45	ASN	2.0
1	D	42	LYS	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 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.

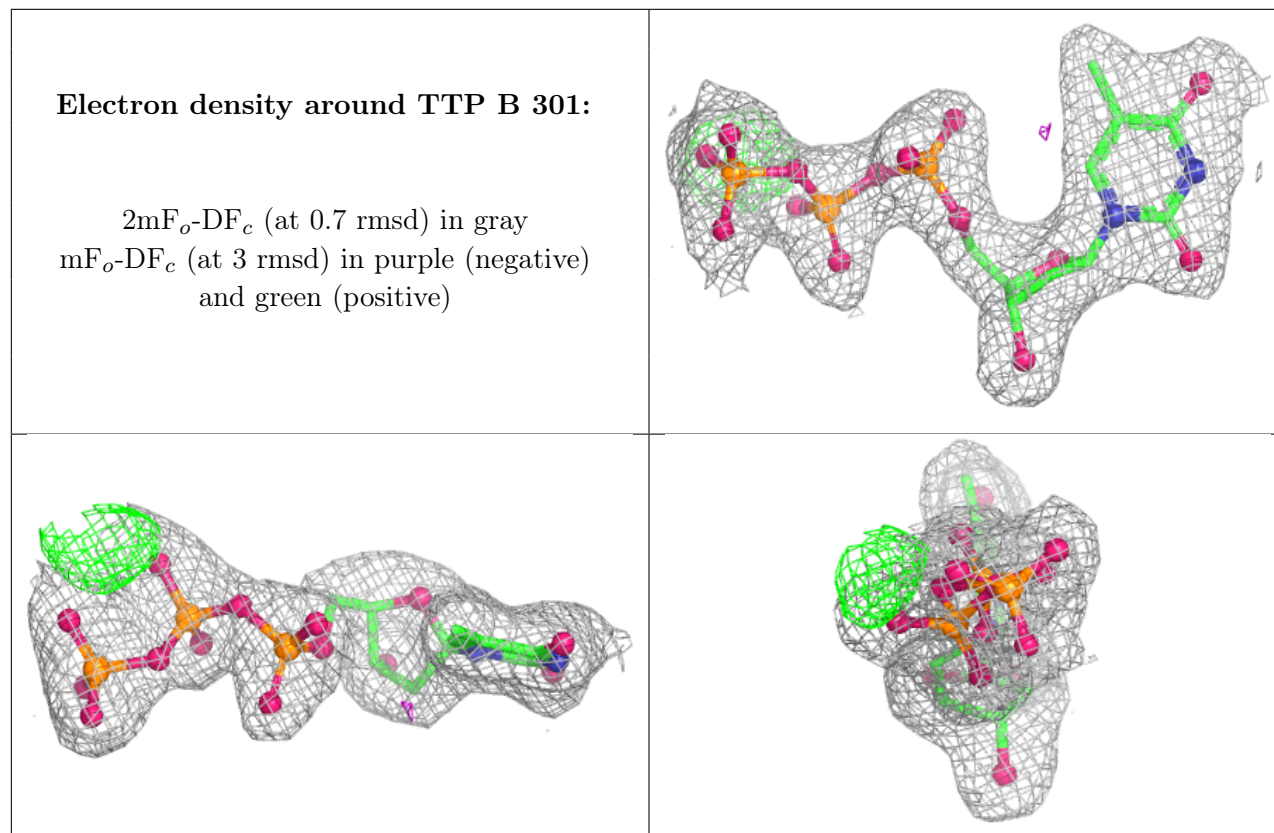
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	MG	D	400	1/1	0.68	0.21	46,46,46,46	0
2	TTP	B	301	29/29	0.97	0.06	16,22,27,27	0
2	TTP	A	300	29/29	0.97	0.06	14,23,28,31	0

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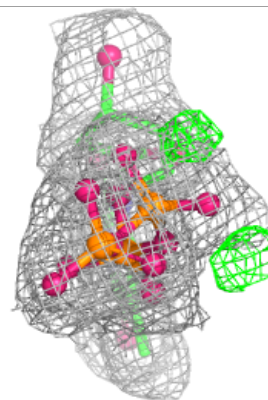
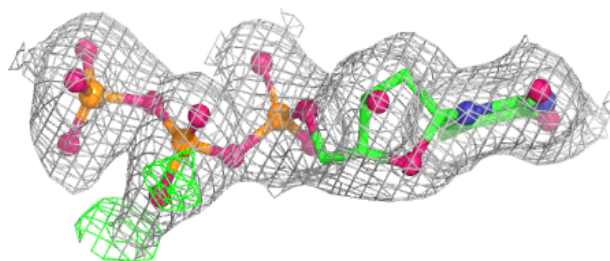
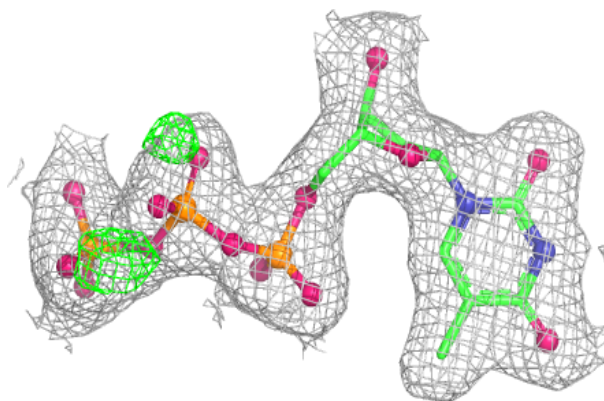
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	TTP	D	303	29/29	0.98	0.05	15,22,26,27	0
2	TTP	C	302	29/29	0.98	0.05	16,22,29,32	0

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.

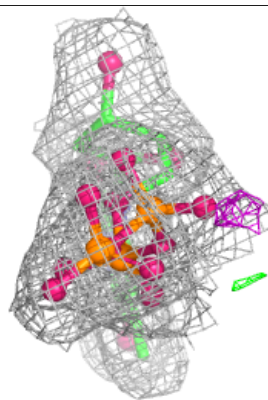
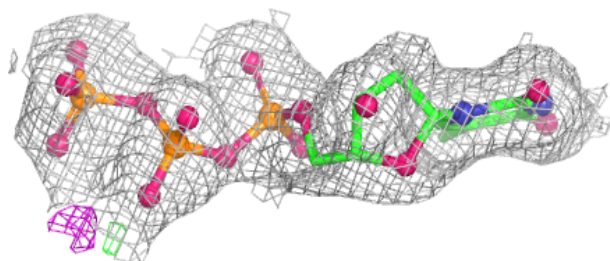
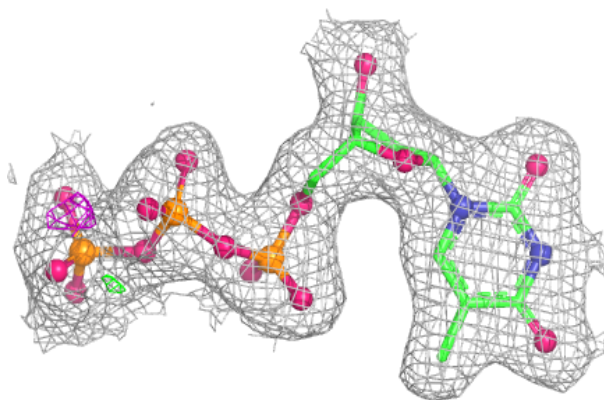


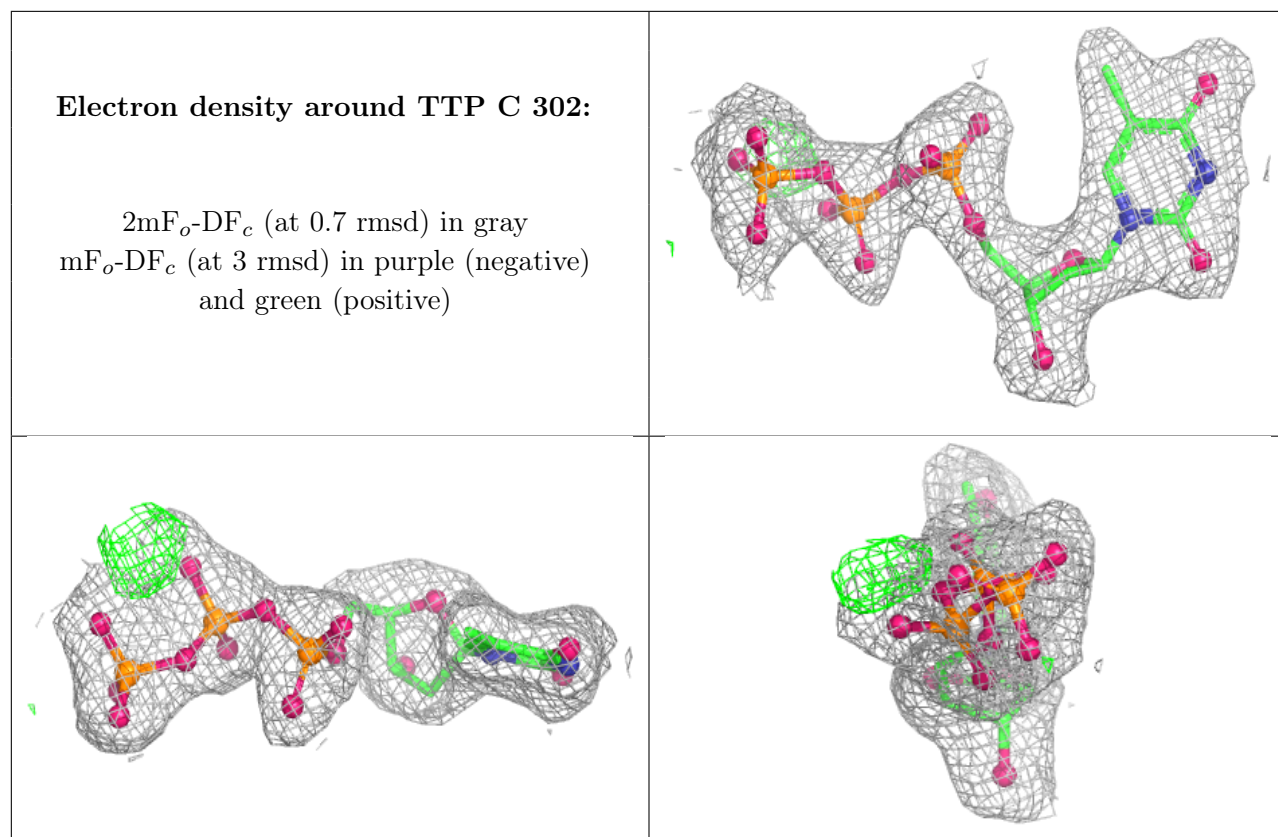
Electron density around TTP A 300:

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

**Electron density around TTP D 303:**

$2mF_o-DF_c$ (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.