



Full wwPDB X-ray Structure Validation Report ⓘ

May 14, 2020 – 08:23 pm BST

PDB ID : 1S0O
Title : Snapshots of replication through an abasic lesion: structural basis for base substitution and frameshift
Authors : Ling, H.; Boudsocq, F.; Woodgate, R.; Yang, W.
Deposited on : 2003-12-31
Resolution : 2.10 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

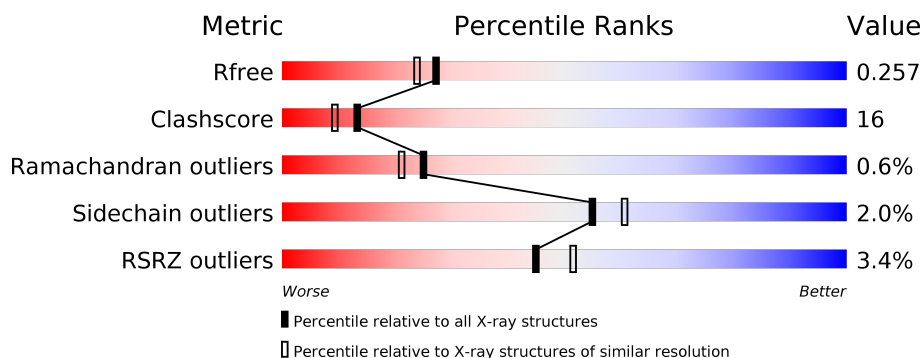
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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	5197 (2.10-2.10)
Clashscore	141614	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (2.10-2.10)
RSRZ outliers	127900	5083 (2.10-2.10)

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	C	13	<div> <div></div> <div>69%31%</div> </div>
1	E	13	<div> <div>8%</div> <div>77%23%</div> </div>
2	D	17	<div> <div>6%</div> <div>29%71%</div> </div>
2	F	17	<div> <div>12%</div> <div>59%29%12%</div> </div>
3	A	352	<div> <div>3%</div> <div>79%17%..</div> </div>
3	B	352	<div> <div>3%</div> <div>70%25%..</div> </div>

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 7124 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 5'-D(*GP*GP*GP*GP*GP*AP*AP*GP*GP*AP*CP*TP*CP)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	C	13	Total	C	N	O	P	0	0	0
			272	128	58	74	12			
1	E	13	Total	C	N	O	P	0	0	0
			272	128	58	74	12			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	17	Total	C	N	O	P	0	0	0
			326	157	52	101	16			
2	F	17	Total	C	N	O	P	0	0	0
			326	157	52	101	16			

- Molecule 3 is a protein called DNA polymerase IV.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	A	341	Total	C	N	O	S	0	0	0
			2743	1760	472	504	7			
3	B	341	Total	C	N	O	S	0	0	0
			2743	1760	472	504	7			

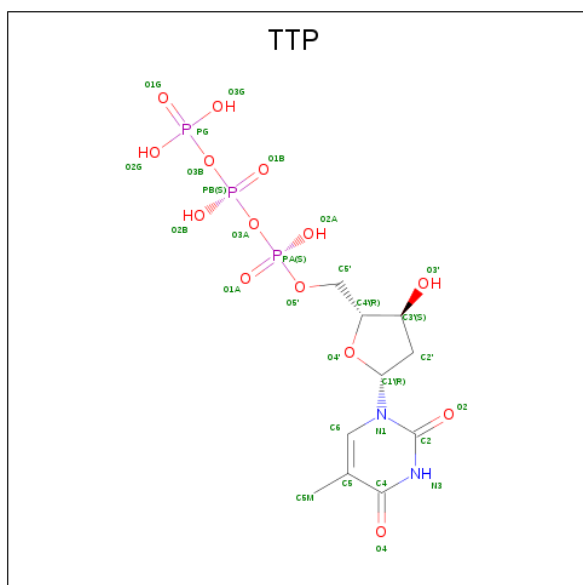
- Molecule 4 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	3	Total	Ca	0	0
			3	3		
4	A	3	Total	Ca	0	0
			3	3		

- Molecule 5 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	B	1	Total Mg 1 1	0	0
5	A	1	Total Mg 1 1	0	0

- Molecule 6 is THYMIDINE-5'-TRIPHOSPHATE (three-letter code: TTP) (formula: $C_{10}H_{17}N_2O_{14}P_3$).



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

- Molecule 7 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	C	18	Total O 18 18	0	0
7	D	23	Total O 23 23	0	0
7	E	32	Total O 32 32	0	0
7	F	25	Total O 25 25	0	0
7	A	149	Total O 149 149	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	B	129	Total 129	O 129	0	0

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.

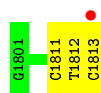
- Molecule 1: 5'-D(*GP*GP*GP*GP*GP*AP*AP*GP*GP*AP*CP*TP*C)-3'

Chain C: 



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

Chain E: 



- Molecule 2: 5'-D(*TP*CP*AP*GP*TP*AP*GP*TP*CP*CP*TP*TP*CP*CP*CP*CP*C)-3',

Chain D: 




- Molecule 2: 5'-D(*TP*CP*AP*GP*TP*AP*GP*TP*CP*CP*TP*TP*CP*CP*CP*CP*C)-3',

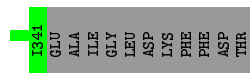
Chain F: 



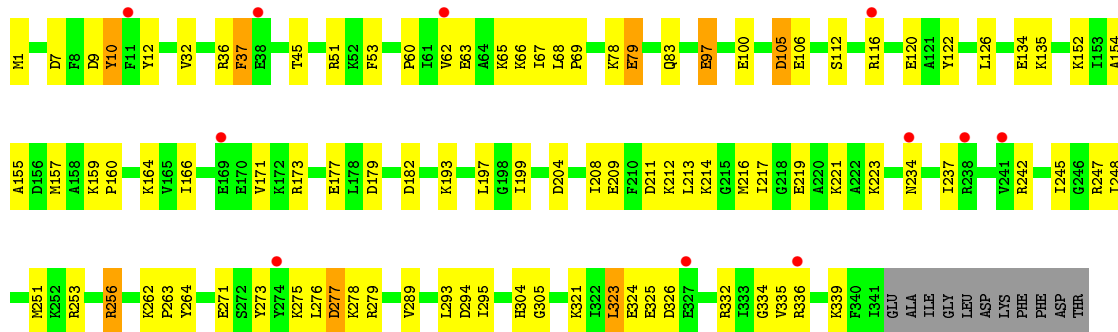
- Molecule 3: DNA polymerase IV

Chain A: 





- Molecule 3: DNA polymerase IV



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	98.30Å 103.05Å 105.91Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.10 29.95 – 2.00	Depositor EDS
% Data completeness (in resolution range)	96.3 (30.00-2.10) 93.0 (29.95-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.37 (at 2.00Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.229 , 0.269 0.219 , 0.257	Depositor DCC
R_{free} test set	2403 reflections (3.43%)	wwPDB-VP
Wilson B-factor (Å ²)	34.9	Xtriage
Anisotropy	0.061	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 43.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.014 for -h,l,k	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7124	wwPDB-VP
Average B, all atoms (Å ²)	41.0	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 34.37 % 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 6.8988e-04. 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: CA, TTP, MG

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	C	0.35	0/307	0.70	0/474
1	E	0.34	0/307	0.74	0/474
2	D	0.34	0/362	0.79	0/554
2	F	0.61	0/362	1.21	5/554 (0.9%)
3	A	0.34	0/2782	0.59	0/3736
3	B	0.34	0/2782	0.60	0/3736
All	All	0.36	0/6902	0.67	5/9528 (0.1%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	1905	DG	OP1-P-O3'	-12.71	77.24	105.20
2	F	1906	DT	C5'-C4'-C3'	6.94	126.60	114.10
2	F	1906	DT	C5'-C4'-O4'	-6.58	96.79	109.30
2	F	1906	DT	O5'-P-OP1	5.76	117.61	110.70
2	F	1906	DT	P-O5'-C5'	-5.38	112.30	120.90

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	C	272	0	146	8	0
1	E	272	0	146	6	0
2	D	326	0	188	27	0
2	F	326	0	188	31	0
3	A	2743	0	2889	62	0
3	B	2743	0	2889	88	1
4	A	3	0	0	0	0
4	B	3	0	0	0	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
6	A	29	0	10	3	0
6	B	29	0	10	6	0
7	A	149	0	0	26	1
7	B	129	0	0	26	1
7	C	18	0	0	7	0
7	D	23	0	0	19	0
7	E	32	0	0	1	0
7	F	25	0	0	7	0
All	All	7124	0	6466	213	2

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 16.

All (213) 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)
6:B:804:TTP:O5'	6:B:804:TTP:C5'	1.64	1.43
3:A:199:ILE:HG12	7:A:952:HOH:O	1.27	1.30
3:A:234:ASN:HB3	7:A:946:HOH:O	1.27	1.26
2:F:1905:DG:H4'	2:F:1906:DT:H5'	1.20	1.18
2:D:1915:DC:H1'	7:D:370:HOH:O	1.48	1.12
1:C:1813:DC:H1'	7:C:349:HOH:O	1.47	1.10
1:C:1801:DG:H5'	3:B:83:GLN:OE1	1.64	0.97
2:F:1905:DG:C4'	2:F:1906:DT:H5'	1.93	0.97
3:A:328:ARG:HB3	7:A:944:HOH:O	1.64	0.96
2:F:1908:DG:H2'	3:B:336:ARG:HH21	1.29	0.95
2:F:1903:DC:O2	2:F:1903:DC:H2'	1.69	0.90
2:F:1902:DT:C4'	7:F:355:HOH:O	2.19	0.89
3:B:116:ARG:HG2	3:B:120:GLU:OE2	1.72	0.89
2:F:1905:DG:H4'	2:F:1906:DT:C5'	2.02	0.89
2:F:1906:DT:H3'	2:F:1907:DA:H5'	1.54	0.88
2:D:1903:DC:O2	2:D:1903:DC:H2'	1.70	0.88

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:234:ASN:CG	7:B:926:HOH:O	2.12	0.87
3:B:9:ASP:HA	7:B:890:HOH:O	1.74	0.85
2:F:1905:DG:N7	7:F:362:HOH:O	2.08	0.84
2:F:1905:DG:C8	7:F:362:HOH:O	2.31	0.84
3:A:290:THR:HB	7:A:944:HOH:O	1.77	0.83
3:A:291:GLU:HG2	7:A:944:HOH:O	1.79	0.81
1:E:1813:DC:H3'	7:B:865:HOH:O	1.79	0.81
3:A:164:LYS:NZ	7:A:951:HOH:O	2.14	0.80
3:B:214:LYS:HE3	3:B:219:GLU:HB2	1.62	0.80
3:A:248:ILE:HD11	3:A:332:ARG:HB3	1.62	0.80
2:D:1909:DT:OP1	2:D:1909:DT:H4'	1.81	0.79
2:F:1905:DG:H2'	3:B:332:ARG:HH22	1.45	0.79
2:F:1908:DG:H2'	3:B:336:ARG:NH2	1.97	0.79
3:A:186:ILE:HG12	7:A:924:HOH:O	1.84	0.78
3:B:212:LYS:HB3	7:B:922:HOH:O	1.83	0.78
3:A:157:MET:HE2	3:A:166:ILE:HD11	1.65	0.77
3:A:45:THR:HG22	7:A:929:HOH:O	1.85	0.76
2:F:1906:DT:H3'	2:F:1907:DA:C5'	2.15	0.76
3:B:7:ASP:CG	7:B:927:HOH:O	2.22	0.76
3:A:10:TYR:HB3	7:A:882:HOH:O	1.87	0.75
3:A:157:MET:CE	3:A:166:ILE:HD11	2.17	0.74
3:B:10:TYR:HA	6:B:804:TTP:O1B	1.88	0.73
3:B:263:PRO:HB2	7:B:907:HOH:O	1.88	0.72
2:D:1914:DC:H2'	7:D:356:HOH:O	1.89	0.71
2:D:1905:DG:O6	7:D:347:HOH:O	2.07	0.71
3:A:224:TYR:HE2	7:A:924:HOH:O	1.73	0.71
1:C:1808:DG:C8	7:C:368:HOH:O	2.43	0.70
3:B:213:LEU:N	7:B:922:HOH:O	2.23	0.70
3:A:77:ARG:HG3	7:A:895:HOH:O	1.91	0.70
2:F:1905:DG:C3'	2:F:1906:DT:H5'	2.18	0.70
3:A:289:VAL:HB	3:A:332:ARG:HB2	1.73	0.69
3:A:248:ILE:CD1	3:A:332:ARG:HB3	2.23	0.68
3:B:273:TYR:HA	3:B:276:LEU:HD12	1.75	0.68
2:D:1914:DC:C2'	7:D:356:HOH:O	2.41	0.68
2:F:1902:DT:O4'	7:F:355:HOH:O	2.12	0.67
3:B:67:ILE:HA	7:B:872:HOH:O	1.95	0.67
2:D:1907:DA:H5''	3:A:242:ARG:HH21	1.60	0.67
1:E:1813:DC:H5'	7:B:806:HOH:O	1.94	0.66
3:B:209:GLU:O	7:B:922:HOH:O	2.14	0.66
3:B:79:GLU:H	3:B:79:GLU:CD	1.97	0.66
2:F:1908:DG:C8	3:B:336:ARG:NH2	2.63	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:154:ALA:HA	3:B:157:MET:HE2	1.77	0.65
3:B:179:ASP:HB3	3:B:182:ASP:OD2	1.97	0.65
3:B:234:ASN:CB	7:B:926:HOH:O	2.44	0.65
3:B:105:ASP:OD1	7:B:865:HOH:O	2.15	0.65
2:D:1912:DT:H2'	7:D:207:HOH:O	1.96	0.65
3:B:208:ILE:HG12	7:B:922:HOH:O	1.97	0.64
6:B:804:TTP:PA	6:B:804:TTP:C5'	2.85	0.64
2:F:1902:DT:H2''	2:F:1903:DC:OP2	1.96	0.64
2:D:1914:DC:C5'	7:D:356:HOH:O	2.45	0.64
3:B:67:ILE:C	3:B:69:PRO:HD3	2.18	0.63
3:B:157:MET:CE	3:B:166:ILE:HD11	2.29	0.63
2:F:1908:DG:H5''	3:B:242:ARG:NH2	2.13	0.62
3:B:97:GLU:CD	3:B:97:GLU:H	2.01	0.62
2:D:1916:DC:H5'	7:D:370:HOH:O	1.99	0.62
3:B:62:VAL:HG23	7:B:848:HOH:O	1.99	0.62
1:C:1807:DA:H2''	7:C:368:HOH:O	1.99	0.61
2:D:1914:DC:H5'	7:D:356:HOH:O	2.00	0.61
3:B:173:ARG:NH1	3:B:177:GLU:OE1	2.31	0.61
3:B:234:ASN:HB3	7:B:926:HOH:O	2.00	0.61
3:A:36:ARG:NH2	3:A:254:ASN:OD1	2.33	0.61
3:A:277:ASP:HB3	7:A:875:HOH:O	2.01	0.61
3:B:157:MET:HE1	3:B:166:ILE:HD11	1.82	0.61
3:A:10:TYR:HA	6:A:803:TTP:O1B	2.01	0.61
3:A:116:ARG:HB2	7:A:923:HOH:O	2.01	0.60
3:A:204:ASP:HB3	7:A:952:HOH:O	2.00	0.60
3:B:105:ASP:OD1	7:B:927:HOH:O	2.16	0.60
3:B:157:MET:HE3	3:B:164:LYS:HD3	1.84	0.60
3:B:304:HIS:HD2	3:B:305:GLY:O	1.85	0.60
3:A:247:ARG:NH1	3:A:249:VAL:HG12	2.17	0.59
3:A:30:VAL:HB	7:A:929:HOH:O	2.01	0.59
3:B:105:ASP:CG	7:B:927:HOH:O	2.40	0.59
1:C:1813:DC:C1'	7:C:349:HOH:O	2.23	0.59
2:D:1914:DC:P	7:D:356:HOH:O	2.60	0.58
3:A:97:GLU:CD	3:A:97:GLU:H	2.06	0.58
2:D:1914:DC:C3'	7:D:356:HOH:O	2.51	0.58
3:B:36:ARG:O	3:B:37:PHE:HB3	2.04	0.58
3:A:51:ARG:NH2	7:A:882:HOH:O	2.36	0.57
3:A:44:ALA:HB3	7:A:929:HOH:O	2.04	0.57
3:A:248:ILE:HD13	3:A:249:VAL:N	2.20	0.56
2:D:1914:DC:H1'	2:D:1915:DC:O5'	2.05	0.56
2:F:1906:DT:OP2	2:F:1907:DA:H3'	2.05	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:1905:DG:H2'	3:B:332:ARG:NH2	2.20	0.56
1:E:1812:DT:H2'	1:E:1813:DC:C6	2.40	0.55
2:F:1902:DT:H4'	7:F:355:HOH:O	1.91	0.55
3:B:245:ILE:HD11	3:B:279:ARG:CZ	2.37	0.54
1:E:1811:DC:H2''	1:E:1812:DT:O5'	2.08	0.54
3:B:248:ILE:HA	3:B:334:GLY:HA3	1.88	0.54
2:D:1908:DG:H2''	2:D:1909:DT:O5'	2.08	0.54
3:B:62:VAL:O	3:B:66:LYS:HG2	2.09	0.53
3:A:36:ARG:NH2	3:A:331:ARG:HG3	2.24	0.53
3:B:289:VAL:HB	3:B:332:ARG:HB2	1.90	0.53
2:D:1913:DT:O3'	7:D:356:HOH:O	2.18	0.53
2:F:1903:DC:C2	3:B:293:LEU:HD11	2.44	0.53
3:B:217:ILE:HD12	3:B:221:LYS:HB3	1.91	0.52
3:B:36:ARG:HD3	7:B:930:HOH:O	2.08	0.52
3:B:155:ALA:HB1	7:B:884:HOH:O	2.09	0.52
3:B:321:LYS:HG2	3:B:325:GLU:OE2	2.09	0.52
2:D:1914:DC:H1'	2:D:1915:DC:C5'	2.39	0.51
3:A:248:ILE:C	3:A:248:ILE:HD13	2.29	0.51
3:A:321:LYS:HG2	7:A:911:HOH:O	2.10	0.51
3:B:335:VAL:HG22	3:B:336:ARG:N	2.24	0.51
3:B:166:ILE:HG22	3:B:171:VAL:HG23	1.92	0.51
3:B:122:TYR:CE2	3:B:126:LEU:HD11	2.45	0.51
2:F:1905:DG:C3'	2:F:1906:DT:C5'	2.89	0.51
3:B:65:LYS:O	3:B:69:PRO:HG3	2.12	0.50
2:D:1903:DC:O2	2:D:1903:DC:C2'	2.49	0.50
3:B:9:ASP:O	3:B:10:TYR:C	2.50	0.50
3:A:1:MET:HE1	7:A:874:HOH:O	2.10	0.50
3:A:262:LYS:HB2	3:A:263:PRO:HD3	1.94	0.49
2:D:1915:DC:H2''	2:D:1916:DC:OP2	2.12	0.49
2:F:1904:DA:H1'	3:B:32:VAL:HG11	1.93	0.49
2:F:1902:DT:C5'	7:F:355:HOH:O	2.53	0.49
3:B:154:ALA:HA	3:B:157:MET:CE	2.42	0.49
2:F:1905:DG:C4'	2:F:1906:DT:C5'	2.76	0.48
3:A:171:VAL:O	3:A:175:ILE:HG13	2.13	0.48
3:B:134:GLU:OE1	3:B:135:LYS:HE2	2.12	0.48
3:A:31:CYS:HB3	3:A:61:ILE:CD1	2.43	0.48
3:B:242:ARG:HG3	3:B:242:ARG:HH11	1.78	0.48
3:B:60:PRO:HD2	3:B:63:GLU:HG3	1.95	0.48
3:B:100:GLU:HB2	3:B:237:ILE:HG23	1.95	0.48
3:B:277:ASP:OD1	3:B:278:LYS:HE2	2.13	0.48
3:A:248:ILE:HD11	3:A:332:ARG:CB	2.37	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:12:TYR:OH	3:B:78:LYS:HE3	2.14	0.48
3:A:183:VAL:HG12	7:A:924:HOH:O	2.13	0.48
3:B:197:LEU:HD11	3:B:216:MET:HG2	1.96	0.48
2:F:1905:DG:C2'	3:B:332:ARG:HH22	2.21	0.48
2:D:1904:DA:H5'	7:A:909:HOH:O	2.14	0.48
7:D:371:HOH:O	3:A:60:PRO:HG3	2.14	0.47
3:B:106:GLU:HG2	7:B:927:HOH:O	2.15	0.47
3:B:294:ASP:OD1	3:B:295:ILE:N	2.46	0.47
3:A:6:VAL:O	3:A:106:GLU:HA	2.15	0.47
2:D:1914:DC:H2'	7:D:259:HOH:O	2.15	0.47
1:C:1808:DG:C5'	7:C:368:HOH:O	2.63	0.47
3:A:9:ASP:O	3:A:10:TYR:C	2.53	0.47
2:D:1916:DC:C5'	7:D:370:HOH:O	2.58	0.47
3:A:252:LYS:HB2	7:A:940:HOH:O	2.15	0.46
3:A:45:THR:HB	6:A:803:TTP:H2'2	1.98	0.46
3:A:36:ARG:HG3	3:A:36:ARG:NH1	2.31	0.46
1:C:1807:DA:C2'	7:C:368:HOH:O	2.60	0.46
6:B:804:TTP:PG	7:B:890:HOH:O	2.73	0.46
2:F:1902:DT:C2'	2:F:1903:DC:OP2	2.64	0.46
1:E:1812:DT:H4'	7:E:170:HOH:O	2.15	0.45
3:A:157:MET:HE1	3:A:166:ILE:HD11	1.96	0.45
3:A:208:ILE:HG12	3:A:209:GLU:N	2.31	0.45
2:F:1903:DC:O2	2:F:1903:DC:C2'	2.48	0.45
3:A:292:ASP:OD1	3:A:328:ARG:HD2	2.18	0.45
3:B:159:LYS:HE2	7:B:827:HOH:O	2.16	0.45
6:A:803:TTP:H3'	6:A:803:TTP:O3A	2.17	0.44
3:B:97:GLU:N	3:B:97:GLU:OE2	2.41	0.44
3:B:157:MET:HE2	3:B:166:ILE:HD11	1.98	0.44
3:A:100:GLU:HG3	3:A:238:ARG:O	2.18	0.44
3:A:292:ASP:O	3:A:293:LEU:HB2	2.17	0.43
3:B:45:THR:HB	6:B:804:TTP:H2'2	1.99	0.43
3:A:157:MET:HE3	3:A:164:LYS:HD3	2.00	0.43
3:B:245:ILE:HD11	3:B:279:ARG:NH2	2.33	0.43
6:B:804:TTP:O2B	6:B:804:TTP:H5'2	2.18	0.43
3:A:256:ARG:NH2	3:A:326:ASP:O	2.52	0.43
3:A:51:ARG:CZ	7:A:882:HOH:O	2.66	0.43
3:B:199:ILE:HG23	3:B:204:ASP:HB2	1.99	0.43
2:F:1902:DT:H5'	7:F:355:HOH:O	2.19	0.43
2:D:1910:DC:H5''	7:D:199:HOH:O	2.17	0.43
3:A:133:LEU:O	3:A:137:LYS:HE3	2.19	0.43
7:D:86:HOH:O	3:A:248:ILE:HG23	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:271:GLU:O	3:B:275:LYS:HG3	2.18	0.43
3:B:68:LEU:N	3:B:69:PRO:HD3	2.33	0.42
3:B:242:ARG:NH1	3:B:242:ARG:HG3	2.34	0.42
3:A:289:VAL:HG21	3:A:332:ARG:HD2	2.01	0.42
3:A:248:ILE:HG23	3:A:248:ILE:O	2.19	0.42
1:E:1813:DC:OP2	3:B:152:LYS:HE3	2.18	0.42
3:B:53:PHE:HB2	3:B:68:LEU:HD11	2.02	0.42
2:D:1915:DC:O3'	7:D:370:HOH:O	2.22	0.42
7:D:86:HOH:O	3:A:248:ILE:HD12	2.18	0.42
3:A:247:ARG:NE	3:A:271:GLU:OE1	2.42	0.42
3:B:159:LYS:HB2	3:B:160:PRO:HA	2.02	0.42
2:D:1914:DC:C1'	2:D:1915:DC:O5'	2.68	0.42
3:A:78:LYS:HB3	7:A:889:HOH:O	2.19	0.42
3:B:208:ILE:CD1	7:B:922:HOH:O	2.67	0.42
1:C:1808:DG:O4'	7:C:368:HOH:O	2.22	0.42
3:B:193:LYS:HG3	7:B:916:HOH:O	2.19	0.42
3:B:323:LEU:HA	3:B:323:LEU:HD12	1.85	0.42
2:D:1914:DC:H4'	2:D:1915:DC:OP1	2.20	0.41
3:A:227:SER:HB2	7:A:921:HOH:O	2.20	0.41
2:D:1914:DC:C4'	7:D:356:HOH:O	2.69	0.41
2:F:1906:DT:O5'	3:B:247:ARG:HD2	2.20	0.41
3:A:208:ILE:HG13	7:A:950:HOH:O	2.21	0.41
3:B:219:GLU:HG3	3:B:223:LYS:HE3	2.03	0.41
3:B:97:GLU:CD	3:B:97:GLU:N	2.72	0.41
3:B:262:LYS:HB2	3:B:263:PRO:HD3	2.03	0.41
3:A:219:GLU:O	3:A:223:LYS:HG3	2.21	0.41
3:B:1:MET:HA	3:B:112:SER:OG	2.21	0.41
3:B:339:LYS:HE3	7:B:856:HOH:O	2.20	0.41
3:B:256:ARG:NH2	3:B:326:ASP:O	2.53	0.41
3:B:251:MET:HA	3:B:264:TYR:CE1	2.56	0.40
2:F:1908:DG:H5''	3:B:242:ARG:CZ	2.51	0.40
3:B:67:ILE:HG23	7:B:872:HOH:O	2.20	0.40
3:A:302:PHE:HZ	3:A:314:GLU:HG2	1.86	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:A:897:HOH:O	7:A:950:HOH:O[3_645]	2.14	0.06
3:B:324:GLU:OE1	7:B:926:HOH:O[4_566]	2.18	0.02

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	339/352 (96%)	325 (96%)	13 (4%)	1 (0%)	41	41
3	B	339/352 (96%)	324 (96%)	12 (4%)	3 (1%)	17	12
All	All	678/704 (96%)	649 (96%)	25 (4%)	4 (1%)	25	21

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	B	277	ASP
3	A	10	TYR
3	B	37	PHE
3	B	10	TYR

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	300/309 (97%)	296 (99%)	4 (1%)	69	75
3	B	300/309 (97%)	292 (97%)	8 (3%)	44	48
All	All	600/618 (97%)	588 (98%)	12 (2%)	55	60

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	A	49	GLU
3	A	105	ASP

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Mol	Chain	Res	Type
3	A	248	ILE
3	A	323	LEU
3	B	51	ARG
3	B	79	GLU
3	B	97	GLU
3	B	105	ASP
3	B	211	ASP
3	B	253	ARG
3	B	256	ARG
3	B	323	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
3	A	188	ASN
3	A	234	ASN
3	B	82	GLN
3	B	285	HIS
3	B	304	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 10 ligands modelled in this entry, 8 are monoatomic - leaving 2 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	803	4	23,30,30	5.99	7 (30%)	29,47,47	2.98	13 (44%)
6	TTP	B	804	4	23,30,30	5.84	6 (26%)	29,47,47	2.81	13 (44%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	TTP	A	803	4	-	2/19/34/34	0/2/2/2
6	TTP	B	804	4	-	2/19/34/34	0/2/2/2

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	803	TTP	C5M-C5	-26.97	0.99	1.51
6	B	804	TTP	C5M-C5	-26.38	1.00	1.51
6	A	803	TTP	PA-O1A	-5.71	1.30	1.50
6	B	804	TTP	PA-O1A	-5.38	1.31	1.50
6	B	804	TTP	O5'-C5'	5.01	1.64	1.44
6	A	803	TTP	O5'-C5'	4.80	1.63	1.44
6	A	803	TTP	O4-C4	3.52	1.33	1.24
6	B	804	TTP	O4-C4	2.90	1.31	1.24
6	B	804	TTP	C4-N3	2.65	1.37	1.33
6	A	803	TTP	C4-N3	2.55	1.37	1.33
6	A	803	TTP	C1'-N1	2.28	1.56	1.49
6	A	803	TTP	C4-C5	2.27	1.46	1.41
6	B	804	TTP	C1'-N1	2.04	1.55	1.49

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	B	804	TTP	C4-N3-C2	7.09	121.13	115.14
6	A	803	TTP	C4-N3-C2	6.89	120.96	115.14
6	A	803	TTP	PB-O3A-PA	-6.44	110.73	132.83
6	A	803	TTP	C2'-C1'-N1	-6.09	100.23	114.27

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	803	TTP	O5'-PA-O1A	-6.08	85.32	109.07
6	B	804	TTP	PB-O3A-PA	-5.76	113.06	132.83
6	B	804	TTP	C2'-C1'-N1	-5.68	101.16	114.27
6	B	804	TTP	O5'-PA-O1A	-5.63	87.08	109.07
6	A	803	TTP	O3G-PG-O3B	4.27	118.97	104.64
6	A	803	TTP	O3B-PG-O1G	-3.89	89.60	111.19
6	B	804	TTP	O3G-PG-O3B	3.43	116.14	104.64
6	A	803	TTP	PB-O3B-PG	3.35	144.31	132.83
6	B	804	TTP	O3B-PG-O1G	-3.31	92.83	111.19
6	B	804	TTP	PB-O3B-PG	2.74	142.22	132.83
6	A	803	TTP	O4'-C1'-C2'	2.59	111.13	106.25
6	B	804	TTP	O2G-PG-O3B	2.48	112.95	104.64
6	B	804	TTP	O4'-C4'-C5'	2.45	117.42	109.37
6	B	804	TTP	O4'-C1'-C2'	2.44	110.86	106.25
6	A	803	TTP	O2G-PG-O3B	2.42	112.74	104.64
6	B	804	TTP	O5'-C5'-C4'	2.39	117.22	108.99
6	A	803	TTP	C6-N1-C1'	2.35	124.52	119.24
6	B	804	TTP	C5-C6-N1	2.31	124.68	122.19
6	A	803	TTP	O3G-PG-O2G	2.16	115.90	107.64
6	A	803	TTP	O5'-C5'-C4'	2.12	116.30	108.99
6	A	803	TTP	O4'-C4'-C5'	2.12	116.34	109.37
6	B	804	TTP	C5M-C5-C6	2.05	123.01	118.68

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	A	803	TTP	PG-O3B-PB-O1B
6	B	804	TTP	PG-O3B-PB-O1B
6	A	803	TTP	C5'-O5'-PA-O1A
6	B	804	TTP	C5'-O5'-PA-O1A

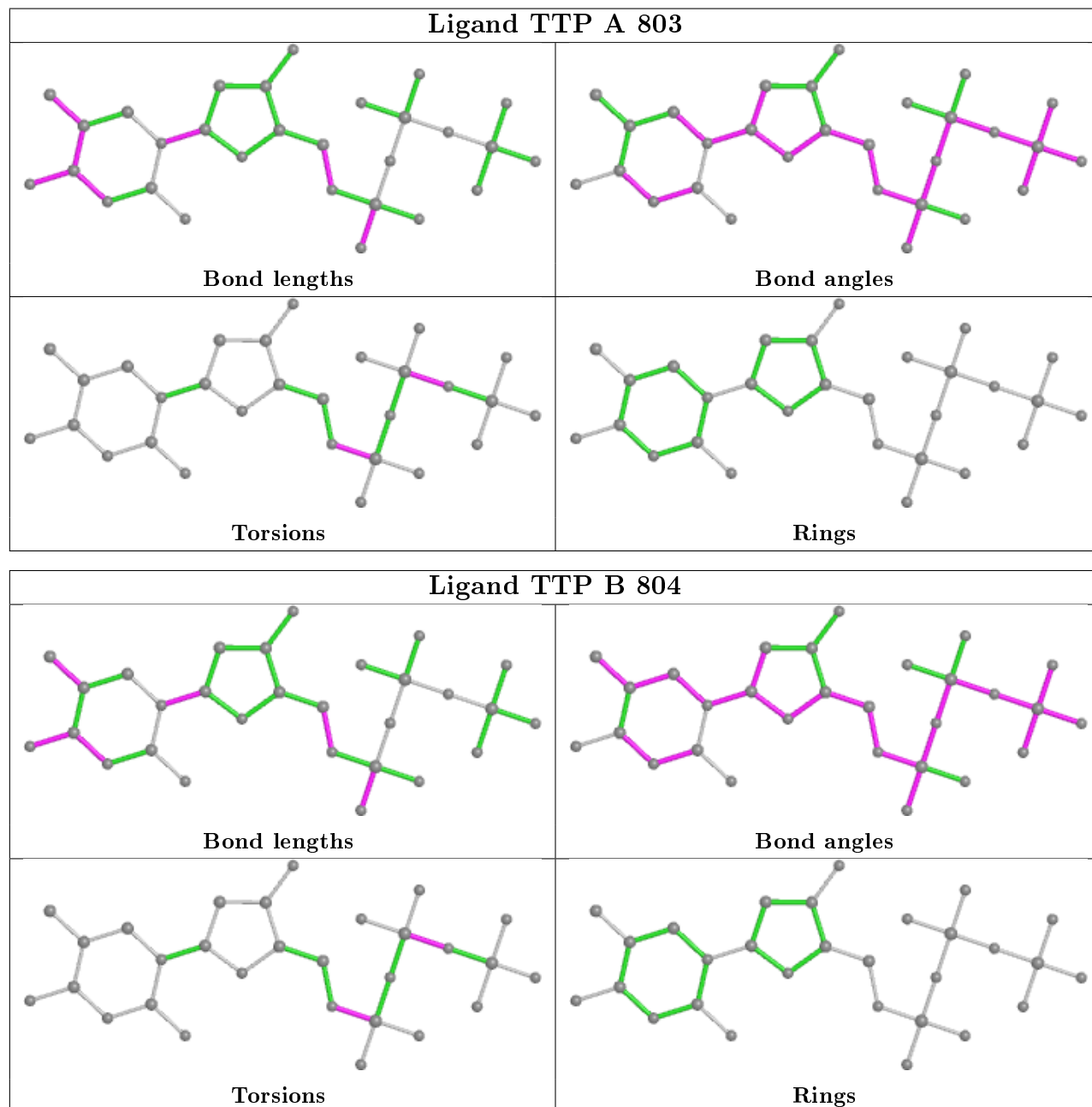
There are no ring outliers.

2 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	803	TTP	3	0
6	B	804	TTP	6	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

There are no such residues in this entry.

5.8 Polymer linkage issues

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	C	13/13 (100%)	0.07	0 100 100	37, 47, 65, 66	0
1	E	13/13 (100%)	0.11	1 (7%) 13 17	33, 38, 51, 73	0
2	D	17/17 (100%)	0.45	1 (5%) 22 27	41, 52, 88, 96	0
2	F	17/17 (100%)	0.43	2 (11%) 4 5	33, 40, 83, 97	0
3	A	341/352 (96%)	0.17	10 (2%) 51 57	21, 36, 54, 68	0
3	B	341/352 (96%)	0.26	11 (3%) 47 54	24, 37, 56, 62	0
All	All	742/764 (97%)	0.22	25 (3%) 45 51	21, 37, 57, 97	0

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	F	1902	DT	3.6
3	B	62	VAL	3.4
3	B	116	ARG	3.3
3	B	274	TYR	3.3
1	E	1813	DC	3.3
3	B	234	ASN	3.1
3	A	238	ARG	3.0
3	A	278	LYS	2.8
3	B	238	ARG	2.7
3	B	241	VAL	2.5
3	A	97	GLU	2.5
3	A	216	MET	2.5
3	A	274	TYR	2.4
3	B	336	ARG	2.3
3	A	169	GLU	2.3
2	F	1903	DC	2.3
3	A	192	GLU	2.3
3	A	252	LYS	2.2
3	B	11	PHE	2.2

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Mol	Chain	Res	Type	RSRZ
3	B	327	GLU	2.1
3	A	253	ARG	2.1
3	B	169	GLU	2.1
2	D	1902	DT	2.1
3	A	318	LEU	2.0
3	B	38	GLU	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](#)

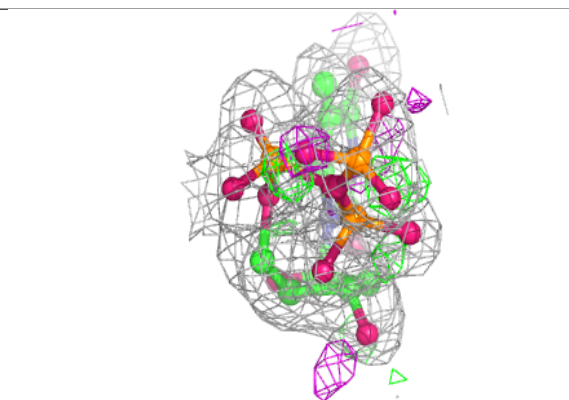
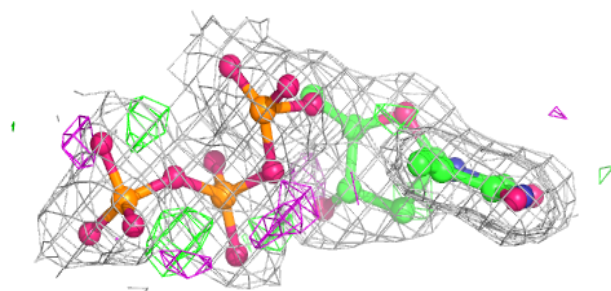
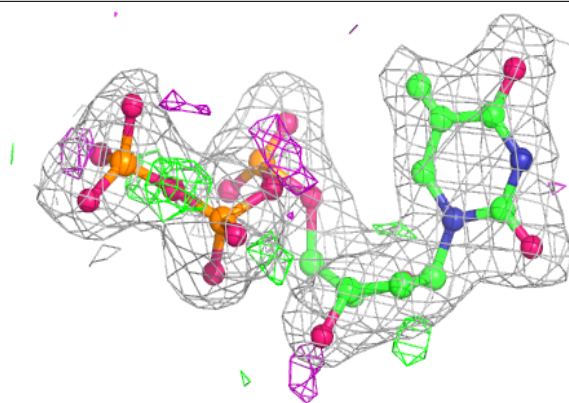
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
5	MG	A	407	1/1	0.38	0.18	45,45,45,45	0
5	MG	B	408	1/1	0.80	0.12	53,53,53,53	0
4	CA	A	406	1/1	0.94	0.10	52,52,52,52	0
4	CA	A	401	1/1	0.96	0.11	36,36,36,36	0
6	TTP	B	804	29/29	0.96	0.16	23,26,38,39	0
6	TTP	A	803	29/29	0.96	0.15	28,32,38,38	0
4	CA	B	403	1/1	0.97	0.08	35,35,35,35	0
4	CA	B	405	1/1	0.97	0.12	41,41,41,41	0
4	CA	B	404	1/1	0.99	0.10	26,26,26,26	0
4	CA	A	402	1/1	0.99	0.15	26,26,26,26	0

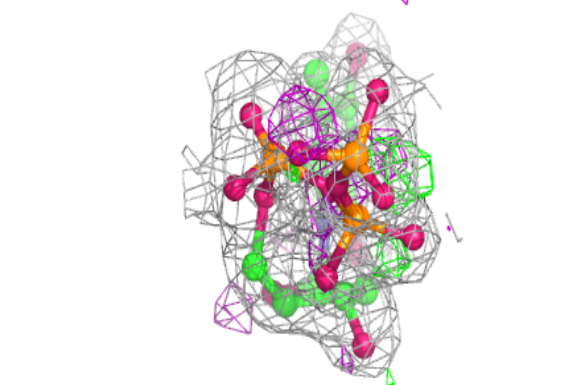
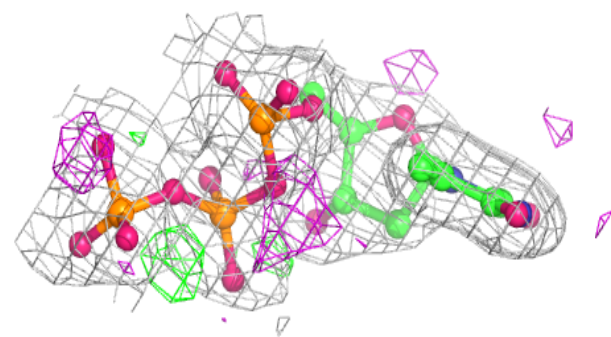
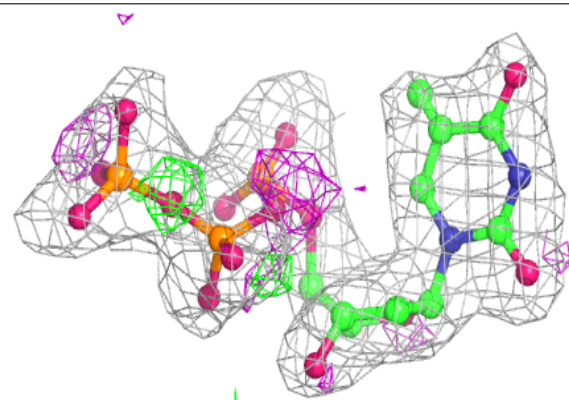
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around TTP B 804:

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

**Electron density around TTP A 803:**

$2mF_o-DF_c$ (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.