



Full wwPDB X-ray Structure Validation Report ⓘ

May 17, 2020 – 01:15 am BST

PDB ID : 4LXT
Title : Crystal structure WlaRD, a sugar 3N-formyl transferase in the presence of dTdp-Qui3N and 5-N-Formyl-THF
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Deposited on : 2013-07-30
Resolution : 1.40 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

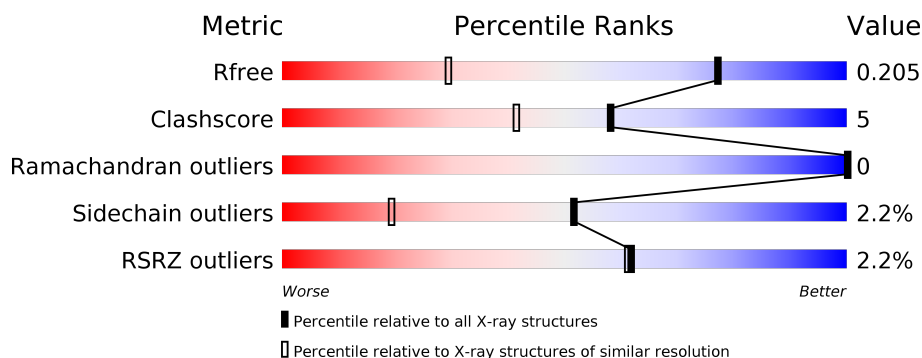
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.40 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	1714 (1.40-1.40)
Clashscore	141614	1812 (1.40-1.40)
Ramachandran outliers	138981	1763 (1.40-1.40)
Sidechain outliers	138945	1762 (1.40-1.40)
RSRZ outliers	127900	1674 (1.40-1.40)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	274	<div> <div></div> <div> <div></div> <div>89%</div> <div>9%</div> </div> </div>
1	B	274	<div> <div>3%</div> <div> <div></div> <div>87%</div> <div>12%</div> </div> </div>

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 5801 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.

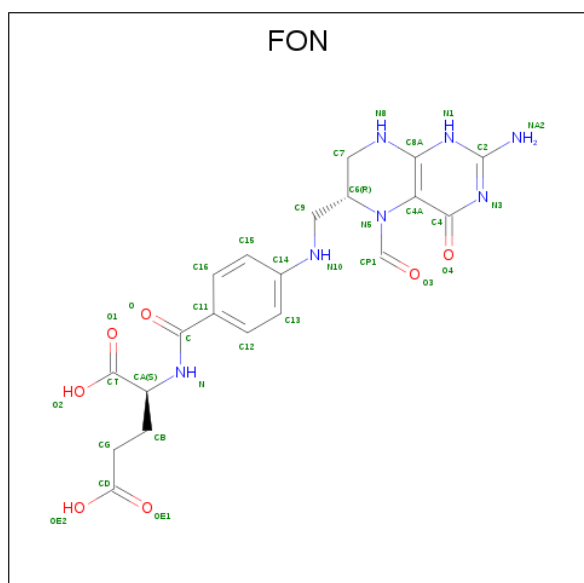
- Molecule 1 is a protein called WlaRD, a sugar 3N-formyl transferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	273	Total	C	N	O	S	0	14	0
			2320	1521	375	421	3			
1	B	274	Total	C	N	O	S	0	10	0
			2313	1514	375	421	3			

There are 4 discrepancies between the modelled and reference sequences:

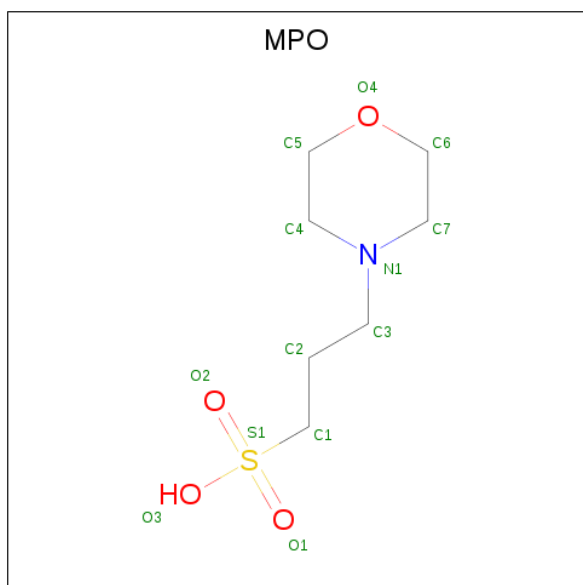
Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	EXPRESSION TAG	UNP A8FMJ3
A	0	HIS	-	EXPRESSION TAG	UNP A8FMJ3
B	-1	GLY	-	EXPRESSION TAG	UNP A8FMJ3
B	0	HIS	-	EXPRESSION TAG	UNP A8FMJ3

- Molecule 2 is N-{[4-({[(6R)-2-amino-5-formyl-4-oxo-1,4,5,6,7,8-hexahydropteridin-6-yl]methyl}amino)phenyl]carbonyl}-L-glutamic acid (three-letter code: FON) (formula: C₂₀H₂₃N₇O₇).



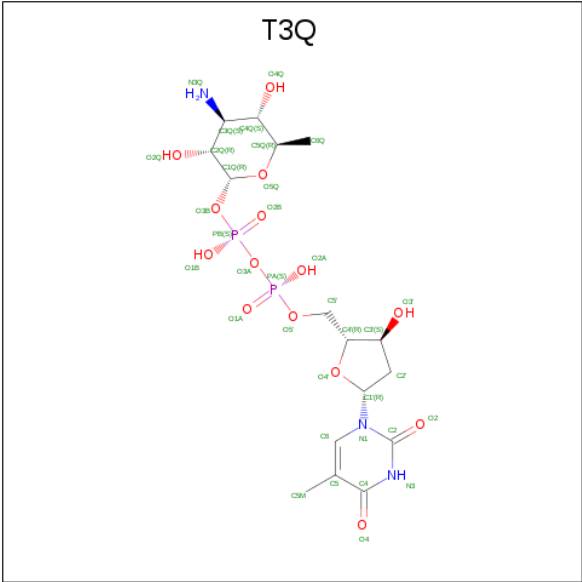
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			34	20	7	7		
2	B	1	Total	C	N	O	0	0
			34	20	7	7		

- Molecule 3 is 3[N-MORPHOLINO]PROPANE SULFONIC ACID (three-letter code: MPO) (formula: C₇H₁₅NO₄S).



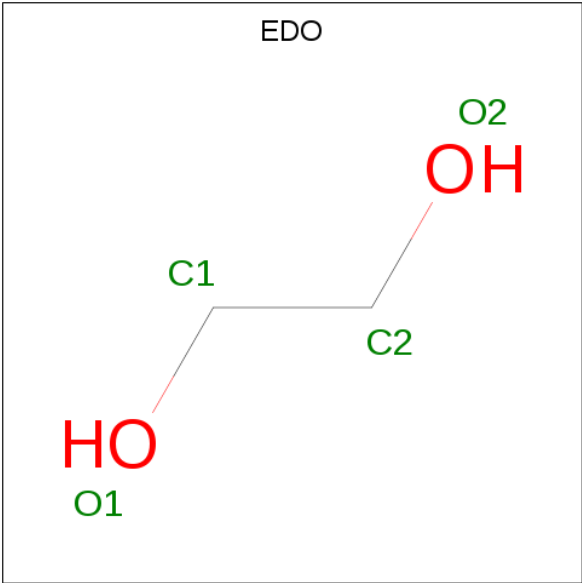
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	S	0	0
			13	7	1	4	1		
3	A	1	Total	C	N	O	S	0	0
			13	7	1	4	1		
3	A	1	Total	C	N	O	S	0	0
			13	7	1	4	1		
3	B	1	Total	C	N	O	S	0	0
			13	7	1	4	1		

- Molecule 4 is [(3R,4S,5S,6R)-4-amino-3,5-dihydroxy-6-methyloxan-2-yl][hydroxy-[[[(2R,3S,5R)-3-hydroxy-5-(5-methyl-2,4-dioxypyrimidin-1-yl)oxolan-2-yl]methoxy]phosphoryl]hydrogen phosphate (three-letter code: T3Q) (formula: C₁₆H₂₇N₃O₁₄P₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	0	0
			35	16	3	14	2		
4	B	1	Total	C	N	O	P	0	0
			35	16	3	14	2		

- Molecule 5 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



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

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

- Molecule 6 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	1	Total 1	Cl 1	0	0
6	A	1	Total 1	Cl 1	0	0

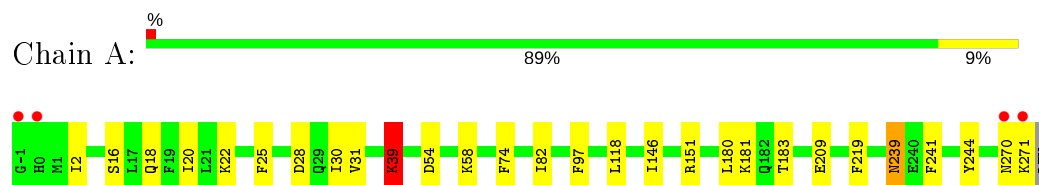
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	526	Total 530	O 530	0	4
7	B	420	Total 422	O 422	0	2

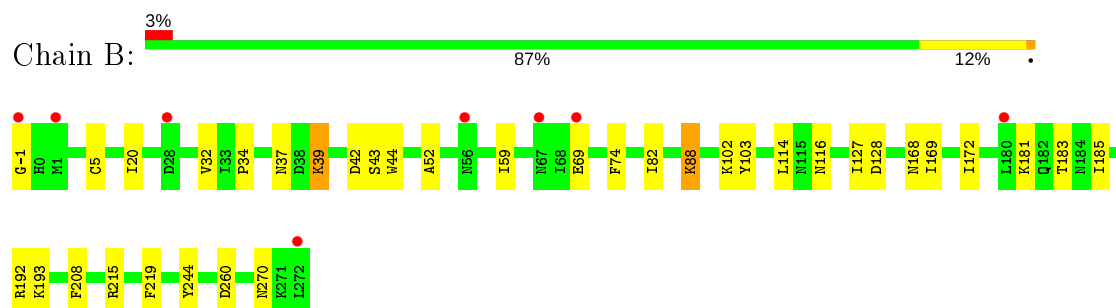
3 Residue-property plots [i](#)

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: WlaRD, a sugar 3N-formyl transferase



- Molecule 1: WlaRD, a sugar 3N-formyl transferase



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	96.78Å 65.15Å 134.72Å 90.00° 110.51° 90.00°	Depositor
Resolution (Å)	30.00 – 1.40 28.69 – 1.40	Depositor EDS
% Data completeness (in resolution range)	92.9 (30.00-1.40) 92.9 (28.69-1.40)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	0.04	Depositor
$\langle I/\sigma(I) \rangle$ ¹	10.26 (at 1.40Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.177 , 0.206 0.177 , 0.205	Depositor DCC
R_{free} test set	7171 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	13.5	Xtriage
Anisotropy	0.020	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 52.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.013 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	5801	wwPDB-VP
Average B, all atoms (Å ²)	18.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.66% 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: MPO, FON, T3Q, EDO, CL

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	0.72	0/2408	1.00	6/3246 (0.2%)
1	B	0.67	1/2392 (0.0%)	1.01	5/3226 (0.2%)
All	All	0.70	1/4800 (0.0%)	1.01	11/6472 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	5	CYS	CB-SG	-5.41	1.73	1.81

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	215	ARG	NE-CZ-NH1	7.43	124.02	120.30
1	B	260	ASP	CB-CG-OD2	-7.12	111.89	118.30
1	A	209	GLU	OE1-CD-OE2	-7.07	114.82	123.30
1	A	39	LYS	CB-CG-CD	-6.43	94.88	111.60
1	A	151	ARG	NE-CZ-NH2	5.59	123.10	120.30
1	B	114	LEU	CB-CG-CD2	-5.58	101.51	111.00
1	A	54	ASP	CB-CG-OD2	-5.49	113.36	118.30
1	B	192	ARG	NE-CZ-NH2	-5.43	117.58	120.30
1	B	128	ASP	CB-CG-OD1	-5.20	113.62	118.30
1	A	28	ASP	CB-CG-OD1	5.01	122.81	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	97	PHE	CB-CG-CD1	5.01	124.31	120.80

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	-1	GLY	Peptide

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2320	0	2395	21	0
1	B	2313	0	2375	21	1
2	A	34	0	21	2	0
2	B	34	0	21	2	0
3	A	39	0	45	8	0
3	B	13	0	15	0	0
4	A	35	0	25	0	0
4	B	35	0	25	0	0
5	A	12	0	18	1	0
5	B	12	0	18	1	0
6	A	1	0	0	0	0
6	B	1	0	0	0	0
7	A	530	0	0	9	1
7	B	422	0	0	4	2
All	All	5801	0	4958	47	3

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 5.

All (47) 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)
7:A:691:HOH:O	1:B:193:LYS:HE3	1.68	0.91

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:32[A]:VAL:HG12	1:B:34:PRO:HD3	1.60	0.81
1:A:82:ILE:H	2:A:301:FON:HN1	1.31	0.79
1:B:82:ILE:H	2:B:301:FON:HN1	1.30	0.76
3:A:304:MPO:H71	7:B:499:HOH:O	1.87	0.74
1:A:39:LYS:HG2	5:A:307:EDO:O1	1.90	0.71
1:B:32[B]:VAL:HG23	1:B:59:ILE:HD12	1.71	0.70
1:A:183[B]:THR:HG21	7:A:571:HOH:O	1.92	0.70
3:A:303:MPO:H42	7:A:610:HOH:O	1.90	0.69
1:A:270:ASN:O	1:A:271:LYS:C	2.30	0.68
3:A:303:MPO:H12	1:B:244:TYR:OH	1.95	0.66
1:B:32[B]:VAL:CG2	1:B:59:ILE:CD1	2.74	0.64
3:A:303:MPO:H71	7:A:453:HOH:O	2.00	0.61
1:A:146:ILE:HD13	1:B:185[B]:ILE:HD12	1.81	0.61
1:B:32[B]:VAL:CG2	1:B:59:ILE:HD12	2.31	0.60
1:B:69:GLU:HG3	1:B:88:LYS:HD3	1.86	0.57
1:A:241:PHE:CE2	3:A:304:MPO:H32	2.41	0.55
1:B:32[A]:VAL:HG12	1:B:34:PRO:CD	2.34	0.54
1:A:16:SER:O	1:A:20[B]:ILE:HG23	2.09	0.53
1:A:118:LEU:HD22	1:B:185[B]:ILE:HD11	1.92	0.51
1:B:44:TRP:HB2	5:B:305:EDO:H11	1.93	0.51
1:B:183[B]:THR:HG21	7:B:481:HOH:O	2.11	0.51
1:A:20[B]:ILE:CD1	1:A:25:PHE:HB2	2.41	0.50
2:B:301:FON:HCP1	2:B:301:FON:O4	2.12	0.49
1:A:180[B]:LEU:HD12	7:A:459:HOH:O	2.11	0.49
1:A:18:GLN:NE2	7:A:834:HOH:O	2.46	0.48
1:B:20[A]:ILE:HD11	1:B:169:ILE:HG12	1.96	0.48
1:B:82:ILE:HD11	1:B:127:ILE:HG21	1.95	0.47
2:A:301:FON:O4	2:A:301:FON:HCP1	2.15	0.46
1:A:244:TYR:OH	3:A:304:MPO:H12	2.16	0.45
1:B:42:ASP:OD2	7:B:767:HOH:O	2.21	0.45
1:B:102:LYS:HB3	1:B:103:TYR:CD2	2.53	0.44
1:A:20[B]:ILE:HD11	1:A:30:ILE:HD11	1.99	0.44
1:B:116:ASN:HB3	1:B:208:PHE:CE2	2.54	0.43
1:A:39:LYS:H	1:A:39:LYS:HG2	1.43	0.42
1:A:22:LYS:HD3	7:A:684:HOH:O	2.20	0.42
3:A:303:MPO:C4	7:A:610:HOH:O	2.60	0.42
1:A:2[A]:ILE:HD12	1:A:31:VAL:HG13	2.01	0.42
3:A:304:MPO:H51	7:B:499:HOH:O	2.20	0.41
1:A:58:LYS:HD3	7:A:744:HOH:O	2.20	0.41
1:B:168:ASN:O	1:B:172[A]:ILE:HG12	2.20	0.41
1:B:52:ALA:HB3	1:B:59:ILE:HD11	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:37:ASN:O	1:B:39:LYS:HD2	2.20	0.41
1:A:20[A]:ILE:HD13	1:A:20[A]:ILE:HG21	1.81	0.40

All (3) 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:B:703:HOH:O	7:B:749:HOH:O[2_556]	2.04	0.16
7:A:741:HOH:O	7:A:824:HOH:O[4_455]	2.07	0.13
1:B:43[B]:SER:OG	7:B:608:HOH:O[2_556]	2.17	0.03

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	285/274 (104%)	276 (97%)	9 (3%)	0	100	100
1	B	283/274 (103%)	273 (96%)	10 (4%)	0	100	100
All	All	568/548 (104%)	549 (97%)	19 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	267/254 (105%)	261 (98%)	6 (2%)	52	19
1	B	265/254 (104%)	259 (98%)	6 (2%)	50	18
All	All	532/508 (105%)	520 (98%)	12 (2%)	52	18

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

Mol	Chain	Res	Type
1	A	39	LYS
1	A	74	PHE
1	A	181	LYS
1	A	219	PHE
1	A	239[A]	ASN
1	A	239[B]	ASN
1	B	39	LYS
1	B	74	PHE
1	B	88	LYS
1	B	181	LYS
1	B	219	PHE
1	B	270	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	80	GLN
1	A	94	ASN
1	B	94	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry

Of 16 ligands modelled in this entry, 2 are monoatomic - leaving 14 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
3	MPO	A	303	-	13,13,13	2.05	1 (7%)	17,17,17	2.66	9 (52%)
5	EDO	A	307	-	3,3,3	0.56	0	2,2,2	0.56	0
5	EDO	A	306	-	3,3,3	0.44	0	2,2,2	0.49	0
5	EDO	A	308	-	3,3,3	0.48	0	2,2,2	0.76	0
3	MPO	B	302	-	13,13,13	2.09	1 (7%)	17,17,17	2.49	7 (41%)
3	MPO	A	302	-	13,13,13	1.81	1 (7%)	17,17,17	2.20	4 (23%)
4	T3Q	A	305	-	31,37,37	1.17	2 (6%)	42,57,57	1.98	7 (16%)
5	EDO	B	307	-	3,3,3	0.48	0	2,2,2	0.69	0
4	T3Q	B	303	-	31,37,37	1.07	2 (6%)	42,57,57	1.62	8 (19%)
5	EDO	B	304	-	3,3,3	0.39	0	2,2,2	0.45	0
2	FON	A	301	-	28,36,36	1.49	3 (10%)	28,50,50	1.43	5 (17%)
5	EDO	B	305	-	3,3,3	0.51	0	2,2,2	0.47	0
2	FON	B	301	-	28,36,36	1.45	2 (7%)	28,50,50	1.45	4 (14%)
3	MPO	A	304	-	13,13,13	2.05	1 (7%)	17,17,17	2.39	7 (41%)

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 Torsions and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	MPO	A	303	-	-	4/7/15/15	0/1/1/1
5	EDO	A	307	-	-	0/1/1/1	-
5	EDO	A	306	-	-	0/1/1/1	-
5	EDO	A	308	-	-	0/1/1/1	-
3	MPO	B	302	-	-	2/7/15/15	0/1/1/1
3	MPO	A	302	-	-	1/7/15/15	0/1/1/1
4	T3Q	A	305	-	-	5/18/53/53	0/3/3/3
5	EDO	B	307	-	-	0/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	T3Q	B	303	-	-	4/18/53/53	0/3/3/3
5	EDO	B	304	-	-	0/1/1/1	-
2	FON	A	301	-	-	2/18/37/37	0/2/3/3
5	EDO	B	305	-	-	1/1/1/1	-
2	FON	B	301	-	-	2/18/37/37	0/2/3/3
3	MPO	A	304	-	-	5/7/15/15	0/1/1/1

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	302	MPO	C1-S1	-7.13	1.67	1.77
3	A	303	MPO	C1-S1	-6.81	1.67	1.77
3	A	304	MPO	C1-S1	-6.78	1.67	1.77
3	A	302	MPO	C1-S1	-6.13	1.68	1.77
2	A	301	FON	CP1-N5	-5.42	1.28	1.35
2	B	301	FON	CP1-N5	-5.10	1.28	1.35
2	B	301	FON	O4-C4	3.99	1.34	1.24
2	A	301	FON	O4-C4	3.60	1.33	1.24
4	B	303	T3Q	C4-N3	3.24	1.38	1.33
4	A	305	T3Q	C6-C5	-2.97	1.31	1.40
4	B	303	T3Q	C6-C5	-2.79	1.32	1.40
4	A	305	T3Q	PA-O2A	-2.43	1.43	1.55
2	A	301	FON	C7-N8	2.01	1.48	1.44

All (51) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	305	T3Q	C4-N3-C2	8.90	122.66	115.14
3	A	302	MPO	C7-N1-C4	6.15	122.67	108.83
3	B	302	MPO	C7-N1-C4	6.08	122.52	108.83
3	A	303	MPO	C7-N1-C4	5.10	120.30	108.83
3	A	302	MPO	C3-N1-C7	4.85	123.64	111.23
4	B	303	T3Q	C4-N3-C2	4.58	119.01	115.14
3	A	304	MPO	C7-N1-C4	4.53	119.04	108.83
3	A	303	MPO	O1-S1-C1	4.47	112.30	106.92
3	A	304	MPO	O1-S1-C1	4.36	112.17	106.92
3	B	302	MPO	O3-S1-C1	4.30	112.72	105.77
3	A	304	MPO	C3-N1-C7	4.02	121.52	111.23
4	B	303	T3Q	O4Q-C4Q-C3Q	-3.94	103.14	110.22
3	A	303	MPO	O3-S1-O1	-3.86	101.83	111.27
4	B	303	T3Q	O4Q-C4Q-C5Q	3.79	118.07	109.67

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	302	MPO	C3-N1-C7	3.78	120.90	111.23
4	A	305	T3Q	O5Q-C5Q-C4Q	-3.77	102.76	109.52
4	A	305	T3Q	O4Q-C4Q-C3Q	-3.69	103.59	110.22
3	A	303	MPO	C3-N1-C7	3.63	120.52	111.23
3	A	303	MPO	C2-C1-S1	-3.47	107.92	113.25
2	A	301	FON	CB-CA-N	-3.33	105.34	110.19
3	A	304	MPO	O3-S1-O1	-3.32	103.16	111.27
2	A	301	FON	C4-N3-C2	3.21	121.04	115.93
3	A	304	MPO	C3-N1-C4	3.21	119.44	111.23
4	B	303	T3Q	O3A-PB-O3B	-3.21	96.02	102.48
3	A	303	MPO	C3-N1-C4	3.11	119.18	111.23
2	B	301	FON	C4-C4A-C8A	2.99	116.78	114.44
4	B	303	T3Q	C5M-C5-C4	-2.98	116.03	121.37
3	B	302	MPO	O1-S1-C1	2.73	110.21	106.92
3	A	302	MPO	C6-C7-N1	2.71	114.21	110.10
3	A	303	MPO	C2-C3-N1	-2.70	107.02	113.84
2	B	301	FON	C4A-N5-C6	-2.67	114.58	119.31
4	A	305	T3Q	C5-C6-N1	2.53	124.92	122.19
3	A	302	MPO	O1-S1-C1	2.47	109.89	106.92
2	B	301	FON	C4-N3-C2	2.42	119.77	115.93
3	B	302	MPO	C5-C4-N1	2.38	113.71	110.10
2	A	301	FON	O3-CP1-N5	-2.33	121.98	125.36
3	A	303	MPO	C6-C7-N1	-2.32	106.59	110.10
4	A	305	T3Q	C2Q-C3Q-C4Q	-2.27	106.12	111.06
3	A	304	MPO	O3-S1-C1	2.27	109.44	105.77
2	B	301	FON	C12-C13-C14	2.27	122.92	120.30
4	B	303	T3Q	C6-N1-C1'	2.25	124.29	119.24
3	A	303	MPO	O3-S1-O2	2.24	116.75	111.27
2	A	301	FON	C15-C16-C11	2.22	123.36	120.78
4	B	303	T3Q	O1B-PB-O2B	2.21	123.15	112.24
3	A	304	MPO	C6-C7-N1	-2.07	106.96	110.10
4	A	305	T3Q	C5M-C5-C6	2.07	123.04	118.68
2	A	301	FON	C4A-C4-N3	-2.04	118.95	123.14
4	B	303	T3Q	O5Q-C5Q-C4Q	-2.04	105.86	109.52
4	A	305	T3Q	C6Q-C5Q-C4Q	2.03	116.83	113.07
3	B	302	MPO	C3-N1-C4	2.03	116.42	111.23
3	B	302	MPO	O3-S1-O1	-2.02	106.34	111.27

There are no chirality outliers.

All (26) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	305	T3Q	C1Q-O3B-PB-O1B
4	B	303	T3Q	C1Q-O3B-PB-O1B
3	A	304	MPO	C2-C1-S1-O1
3	A	304	MPO	C2-C1-S1-O3
3	A	304	MPO	C2-C3-N1-C7
3	A	303	MPO	C2-C3-N1-C7
3	A	302	MPO	C2-C3-N1-C7
3	B	302	MPO	C2-C3-N1-C7
2	A	301	FON	C15-C14-N10-C9
3	A	303	MPO	C2-C1-S1-O3
2	A	301	FON	C13-C14-N10-C9
3	A	304	MPO	C1-C2-C3-N1
2	B	301	FON	C15-C14-N10-C9
5	B	305	EDO	O1-C1-C2-O2
2	B	301	FON	C13-C14-N10-C9
3	A	303	MPO	C2-C1-S1-O1
3	A	303	MPO	C2-C1-S1-O2
3	A	304	MPO	C2-C1-S1-O2
4	A	305	T3Q	PB-O3A-PA-O5'
4	B	303	T3Q	PB-O3A-PA-O5'
4	B	303	T3Q	C1Q-O3B-PB-O3A
4	A	305	T3Q	C1Q-O3B-PB-O2B
4	B	303	T3Q	C1Q-O3B-PB-O2B
4	A	305	T3Q	C1Q-O3B-PB-O3A
3	B	302	MPO	C2-C3-N1-C4
4	A	305	T3Q	PB-O3A-PA-O1A

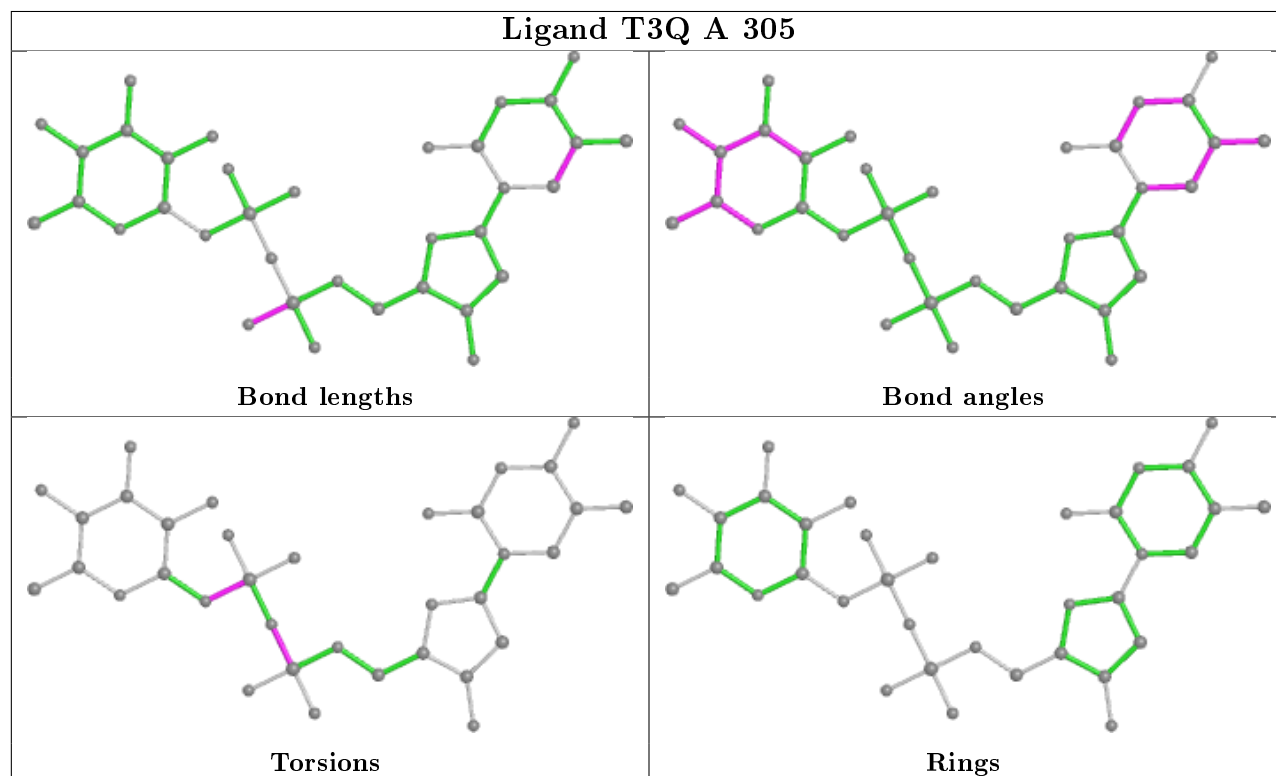
There are no ring outliers.

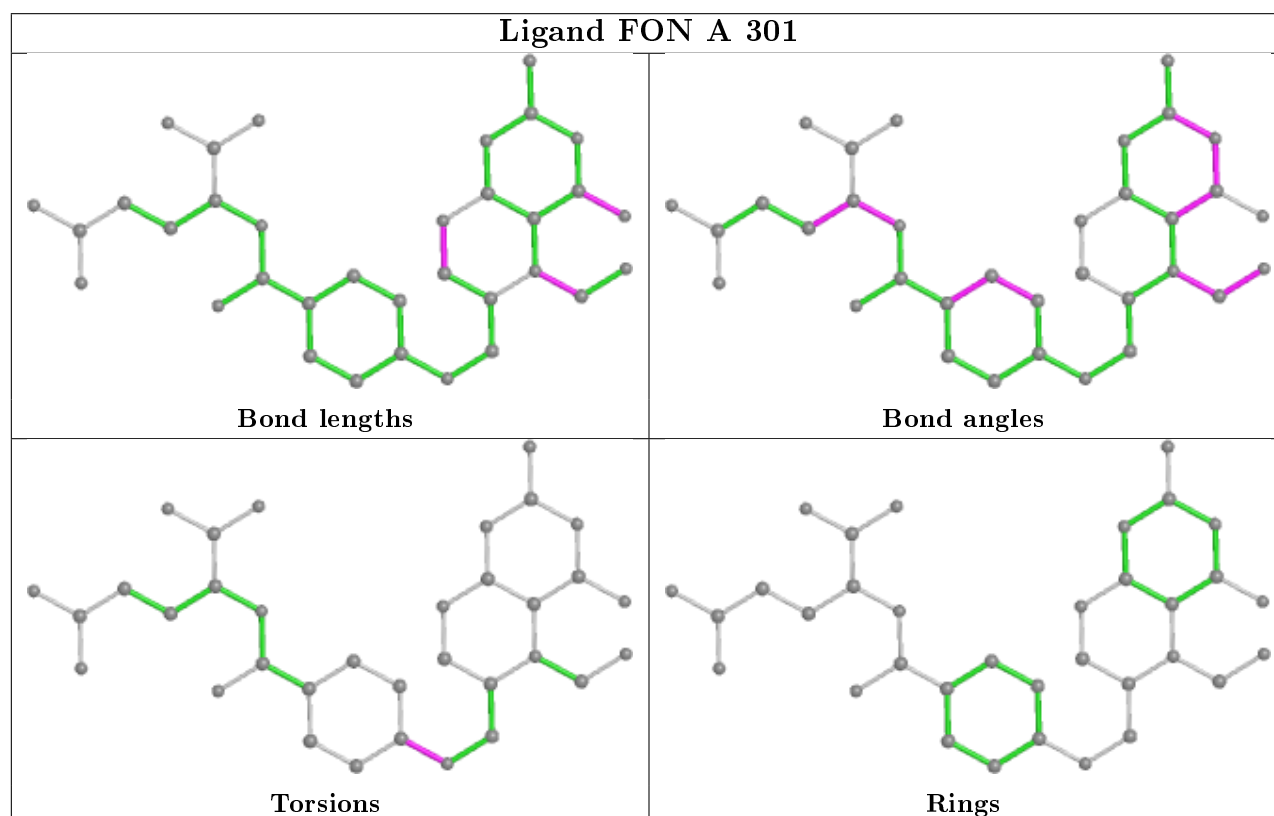
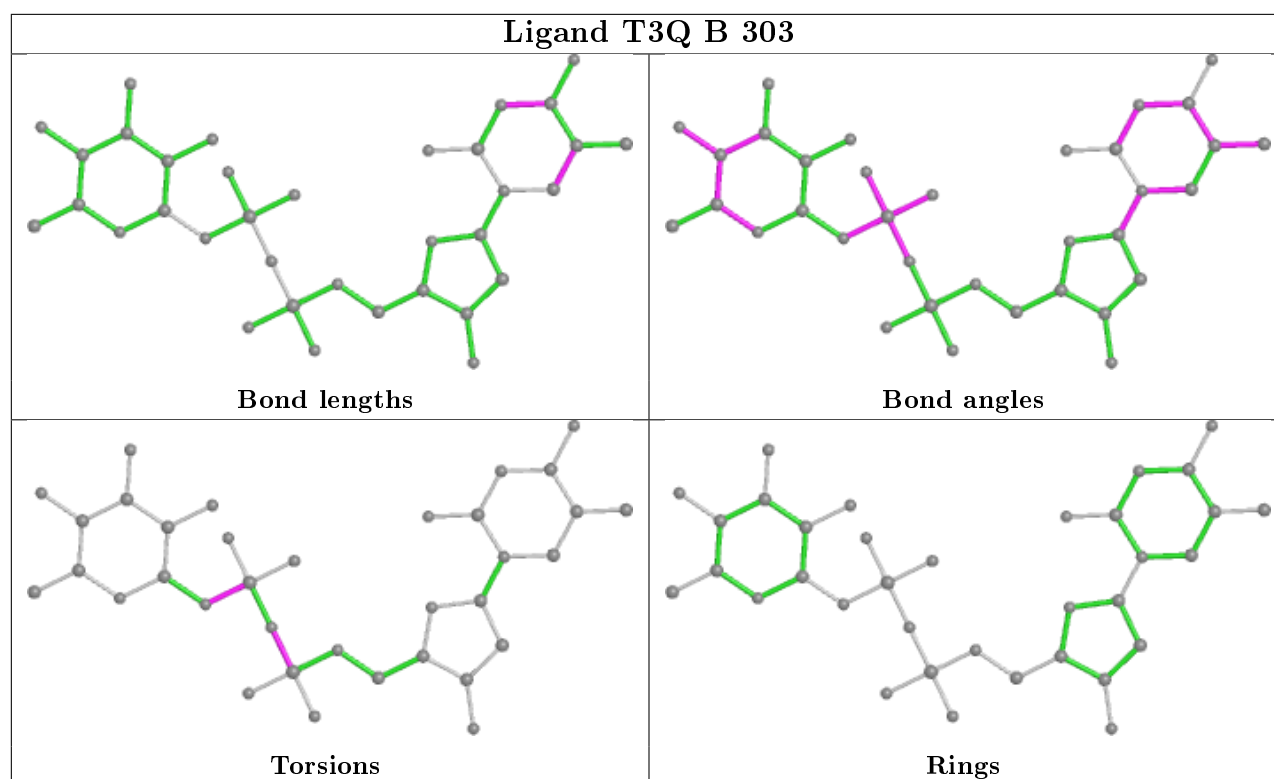
6 monomers are involved in 14 short contacts:

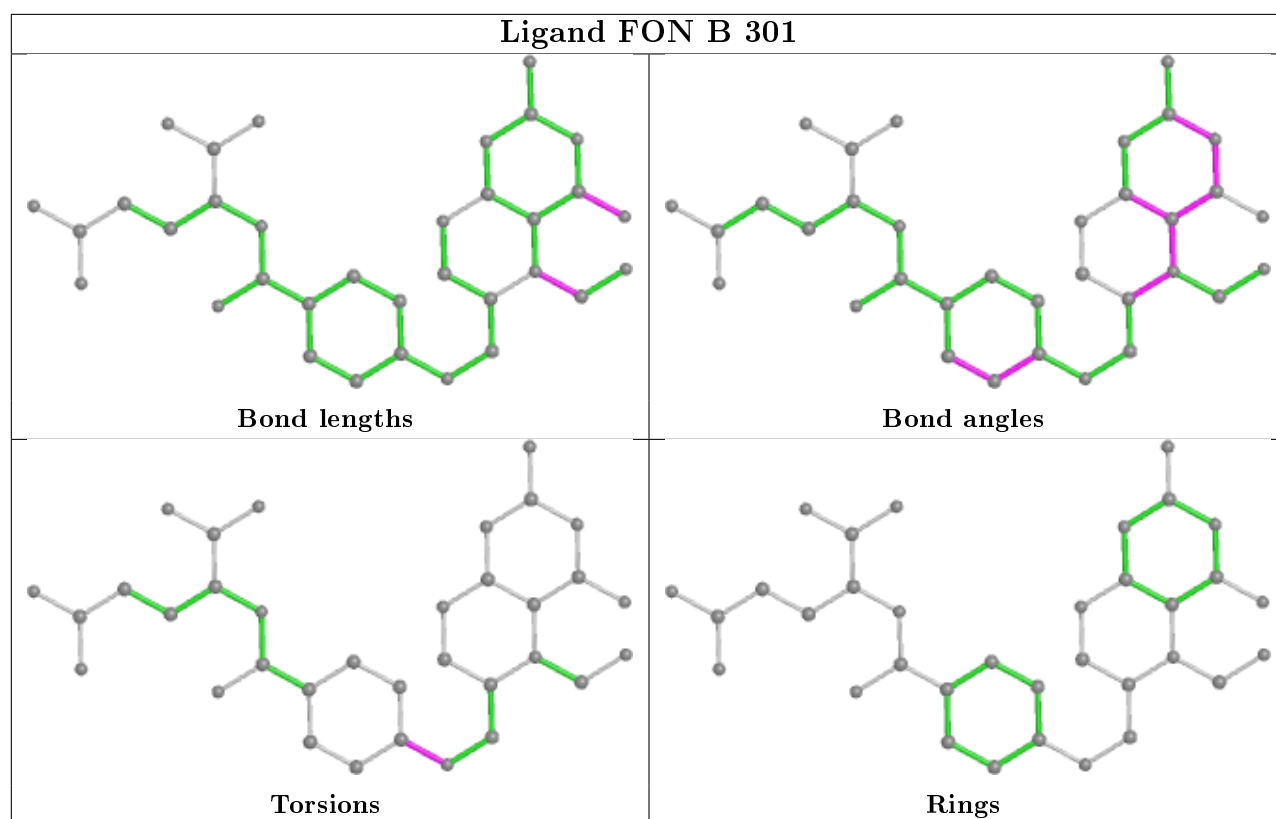
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	303	MPO	4	0
5	A	307	EDO	1	0
2	A	301	FON	2	0
5	B	305	EDO	1	0
2	B	301	FON	2	0
3	A	304	MPO	4	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less 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	273/274 (99%)	-0.40	4 (1%) 73 72	9, 13, 24, 45	0
1	B	274/274 (100%)	-0.09	8 (2%) 51 50	9, 16, 32, 51	0
All	All	547/548 (99%)	-0.24	12 (2%) 62 61	9, 14, 29, 51	0

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	-1	GLY	6.0
1	B	-1	GLY	5.9
1	A	271	LYS	3.1
1	A	270	ASN	3.0
1	B	56[A]	ASN	2.7
1	B	180	LEU	2.4
1	B	67	ASN	2.3
1	B	1	MET	2.2
1	B	28	ASP	2.2
1	A	0	HIS	2.1
1	B	272	LEU	2.1
1	B	69	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

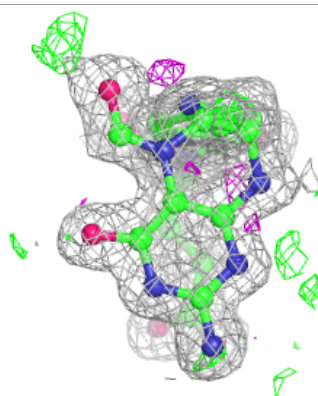
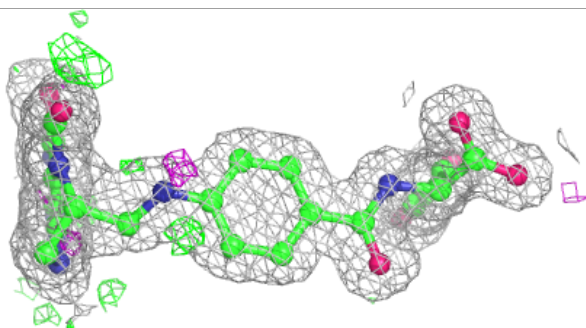
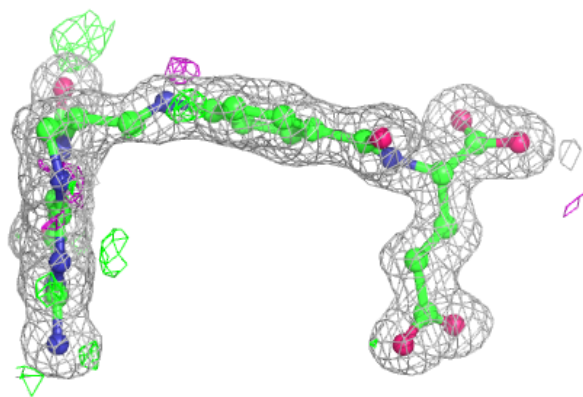
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	MPO	A	303	13/13	0.86	0.20	20,40,56,57	0
3	MPO	A	304	13/13	0.87	0.24	22,43,64,65	0
5	EDO	B	305	4/4	0.93	0.14	26,31,35,36	0
2	FON	B	301	34/34	0.94	0.08	14,16,26,32	0
2	FON	A	301	34/34	0.95	0.08	9,13,20,29	0
5	EDO	A	308	4/4	0.96	0.07	19,21,30,32	0
3	MPO	A	302	13/13	0.96	0.12	19,22,26,27	0
3	MPO	B	302	13/13	0.97	0.10	19,23,29,31	0
5	EDO	A	306	4/4	0.97	0.07	20,20,21,26	0
5	EDO	B	304	4/4	0.98	0.05	16,21,21,25	0
5	EDO	A	307	4/4	0.98	0.08	16,19,21,28	0
4	T3Q	A	305	35/35	0.98	0.06	8,11,14,17	0
5	EDO	B	307	4/4	0.98	0.06	16,22,31,31	0
4	T3Q	B	303	35/35	0.98	0.05	10,12,17,20	0
6	CL	A	309	1/1	0.99	0.03	16,16,16,16	0
6	CL	B	306	1/1	0.99	0.04	21,21,21,21	0

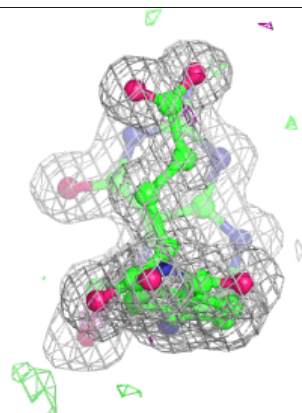
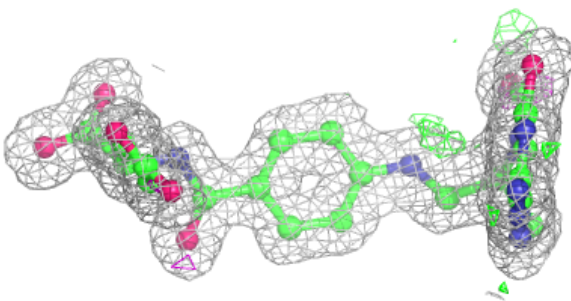
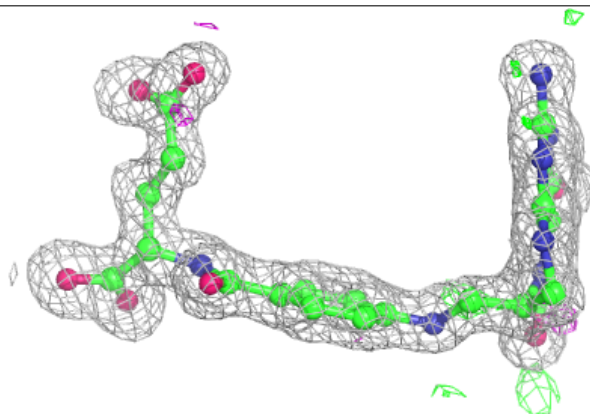
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around FON B 301:

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

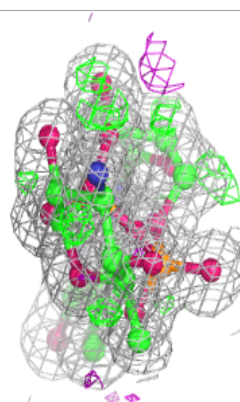
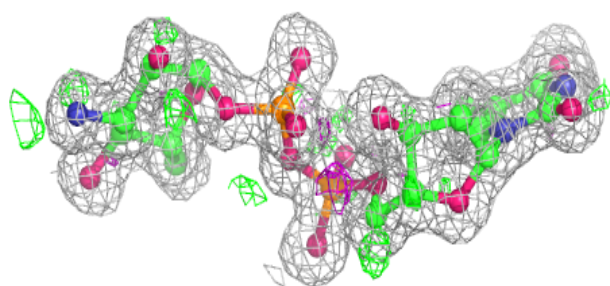
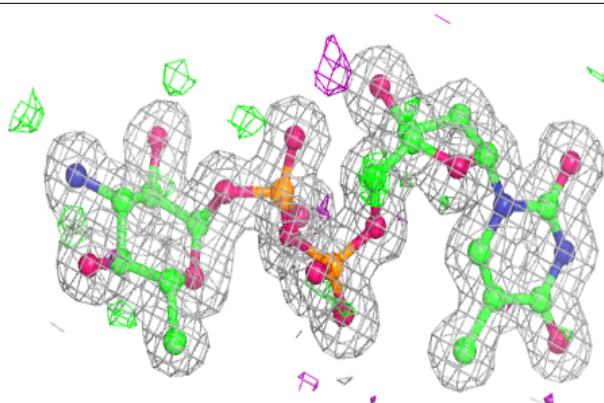
**Electron density around FON A 301:**

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

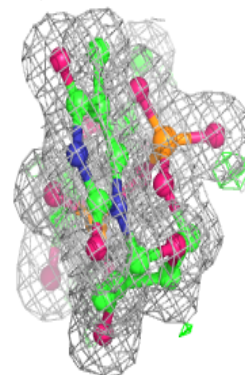
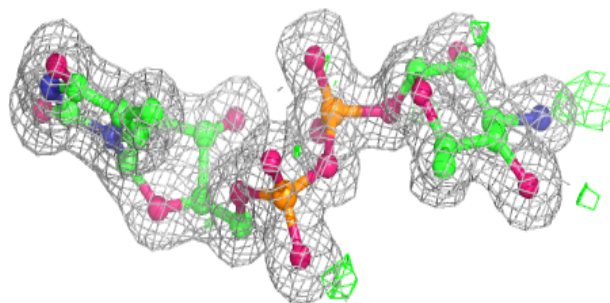
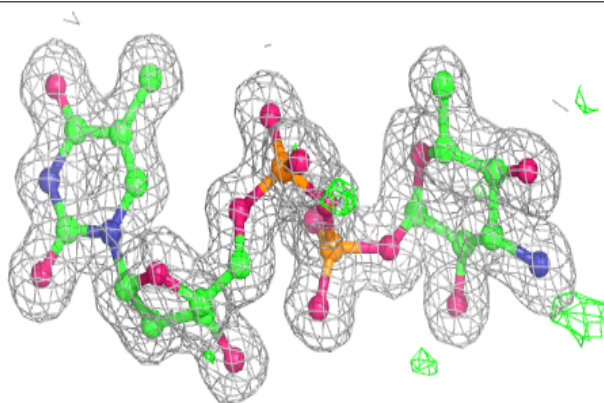


Electron density around T3Q A 305:

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

**Electron density around T3Q B 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.