

Full wwPDB X-ray Structure Validation Report (i)

Sep 11, 2023 – 03:29 PM EDT

PDB ID : 4LY3

Title: Crystal structure of WlaRD, a sugar 3N-formyl transferase in the presence of

dTPD-Qui3N, dTDP-Qui3NFo, and THF

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Deposited on : 2013-07-30

Resolution : 1.90 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 (i) 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 (1)) 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 : 2.35.1

buster-report : 1.1.7 (2018)

Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)

 $Refmac \quad : \quad 5.8.0158$

CCP4 : 7.0.044 (Gargrove)

Ideal geometry (proteins) : Engh & Huber (2001) Ideal geometry (DNA, RNA) : Parkinson et al. (1996)

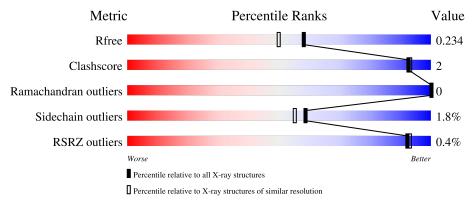
Validation Pipeline (wwPDB-VP) : 2.35.1

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: X- $RAY\ DIFFRACTION$

The reported resolution of this entry is 1.90 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	Similar resolution
Metric	$(\# \mathrm{Entries})$	$(\# ext{Entries}, ext{ resolution range}(ext{Å}))$
R_{free}	130704	6207 (1.90-1.90)
Clashscore	141614	6847 (1.90-1.90)
Ramachandran outliers	138981	6760 (1.90-1.90)
Sidechain outliers	138945	6760 (1.90-1.90)
RSRZ outliers	127900	6082 (1.90-1.90)

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 <=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	95%						
1	В	274	92%	7% •					



2 Entry composition (i)

There are 6 unique types of molecules in this entry. The entry contains 5254 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 formyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	271	10001	C 1462	- '	O 413	S 3	0	2	0
1	В	271	Total 2237	C 1457		O 411	S 3	0	1	0

There are 4 discrepancies between the modelled and reference sequences:

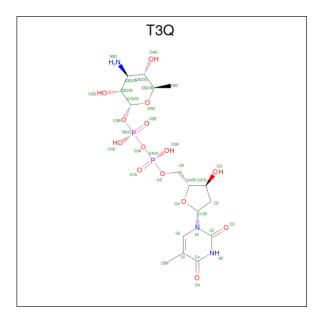
Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	expression tag	UNP A8FMJ3
A	0	HIS	-	expression tag	UNP A8FMJ3
В	-1	GLY	-	expression tag	UNP A8FMJ3
В	0	HIS	-	expression tag	UNP A8FMJ3

• Molecule 2 is N-[4-({[(6R)-2-amino-4-oxo-3,4,5,6,7,8-hexahydropteridin-6-yl]methyl}amino) benzoyl]-L-glutamic acid (three-letter code: 1YJ) (formula: C₁₉H₂₃N₇O₆).



Mol	Chain	Residues					ZeroOcc	AltConf
2	A	1	Total 32		N 7		0	0
2	В	1	Total 32		N 7		0	0

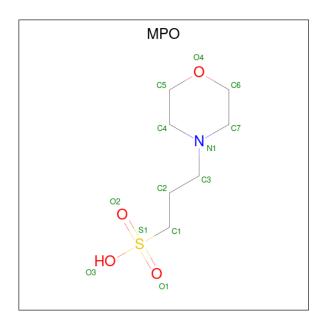
• Molecule 3 is $[(3R,4S,5S,6R)-4-amino-3,5-dihydroxy-6-methyloxan-2-yl][hydroxy-[[(2R,3S,5R)-3-hydroxy-5-(5-methyl-2,4-dioxopyrimidin-1-yl)oxolan-2-yl]methoxy]phosphoryl] hydrogen phosphate (three-letter code: T3Q) (formula: <math>C_{16}H_{27}N_3O_{14}P_2$).



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

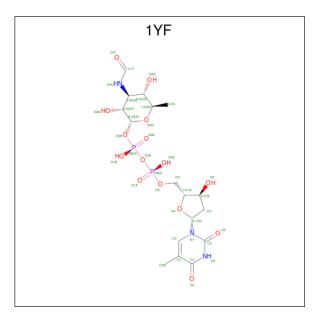
• Molecule 4 is 3[N-MORPHOLINO]PROPANE SULFONIC ACID (three-letter code: MPO) (formula: $C_7H_{15}NO_4S$).





Mol	Chain	Residues	Atoms					ZeroOcc	AltConf			
4	Λ	Λ 1	Total	С	N	О	S	0	0			
4 A	1	13	7	1	4	1	0	0				
1	D	D	D	D	1	Total	С	N	О	S	0	0
4	Б	1	13	7	1	4	1	0				

• Molecule 5 is [(2R,3R,4S,5S,6R)-4-formamido-6-methyl-3,5-bis(oxidanyl)oxan-2-yl] $[[(2R,3S,5R)-5-[5-methyl-2,4-bis(oxidanylidene)pyrimidin-1-yl]-3-oxidanyl-oxolan-2-yl]methoxy-oxid anyl-phosphoryl] hydrogen phosphate (three-letter code: 1YF) (formula: <math>C_{17}H_{27}N_3O_{15}P_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	В	1	Total 37	C 17			P 2	0	0



• Molecule 6 is water.

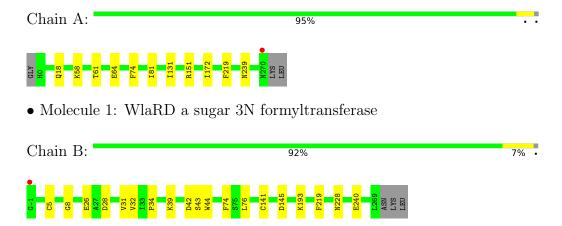
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	308	Total O 308 308	0	0
6	В	301	Total O 301 301	0	0



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: WlaRD a sugar 3N formyltransferase





4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants	96.00Å 65.55Å 134.50Å	Donogitor
a, b, c, α , β , γ	90.00° 109.83° 90.00°	Depositor
Resolution (Å)	50.00 - 1.90	Depositor
Resolution (A)	63.26 - 1.89	EDS
% Data completeness	91.9 (50.00-1.90)	Depositor
(in resolution range)	92.0 (63.26-1.89)	EDS
R_{merge}	0.05	Depositor
R_{sym}	0.05	Depositor
$< I/\sigma(I) > 1$	1.65 (at 1.88Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
D.D.	0.178 , 0.226	Depositor
R, R_{free}	0.187 , 0.234	DCC
R_{free} test set	2940 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	16.0	Xtriage
Anisotropy	0.071	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.34 , 47.5	EDS
L-test for twinning ²	$< L > = 0.48, < L^2> = 0.31$	Xtriage
Estimated twinning fraction	0.023 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5254	wwPDB-VP
Average B, all atoms (Å ²)	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 4.54% of the height of the origin peak. No significant pseudotranslation is detected.

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



¹Intensities estimated from amplitudes.

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: 1YJ, T3Q, 1YF, MPO

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		
IVIOI		RMSZ	# Z > 5	RMSZ	# Z > 5	
1	A	0.65	0/2298	0.93	1/3100 (0.0%)	
1	В	0.66	0/2286	0.94	$2/3083 \ (0.1\%)$	
All	All	0.66	0/4584	0.94	3/6183 (0.0%)	

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$Observed(^o)$	$\mathbf{Ideal}(^{o})$
1	В	28	ASP	CB-CG-OD2	5.88	123.59	118.30
1	A	151	ARG	NE-CZ-NH2	5.01	122.81	120.30
1	В	28	ASP	CB-CG-OD1	-5.01	113.79	118.30

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	2246	0	2266	3	0
1	В	2237	0	2257	10	0
2	A	32	0	21	0	0
2	В	32	0	21	0	0
3	A	35	0	25	0	0
4	A	13	0	15	0	0



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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	В	13	0	15	1	0
5	В	37	0	25	1	0
6	A	308	0	0	1	0
6	В	301	0	0	6	2
All	All	5254	0	4645	14	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 2.

All (14) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic	Clash
Atom-1	Atom-2	${f distance} \; ({ m \AA})$	overlap (Å)
5:B:302:1YF:H2	6:B:418:HOH:O	1.74	0.87
1:B:42:ASP:OD2	6:B:642:HOH:O	1.95	0.83
1:B:141:CYS:SG	6:B:490:HOH:O	2.48	0.72
1:B:240:GLU:OE1	6:B:622:HOH:O	2.10	0.68
1:A:18:GLN:NE2	6:A:638:HOH:O	2.42	0.51
1:B:26:GLU:HB3	6:B:569:HOH:O	2.10	0.50
1:B:43[B]:SER:OG	1:B:44:TRP:N	2.44	0.50
1:B:8:GLY:HA3	1:B:76:LEU:O	2.13	0.49
1:B:228:ASN:OD1	4:B:303:MPO:H12	2.18	0.43
1:A:81:ILE:HD12	1:A:131:ILE:HG12	2.01	0.42
1:B:32:VAL:HG23	1:B:34:PRO:HD3	2.02	0.41
1:A:61:THR:OG1	1:A:64:GLU:HG3	2.20	0.41
1:B:5:CYS:HA	1:B:31:VAL:O	2.21	0.41
1:B:145:ASP:HB3	6:B:595:HOH:O	2.21	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	$\begin{array}{c} {\rm Interatomic} \\ {\rm distance} \ ({\rm \AA}) \end{array}$	Clash overlap (Å)
6:B:541:HOH:O	6:B:650:HOH:O[3_445]	2.15	0.05
6:B:416:HOH:O	6:B:650:HOH:O[3_445]	2.16	0.04



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	Perce	ntiles
1	A	271/274~(99%)	258 (95%)	13 (5%)	0	100	100
1	В	$270/274\ (98\%)$	260 (96%)	10 (4%)	0	100	100
All	All	541/548 (99%)	518 (96%)	23 (4%)	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	$254/254 \ (100\%)$	249 (98%)	5 (2%)	55 51		
1	В	252/254~(99%)	248 (98%)	4 (2%)	62 60		
All	All	506/508 (100%)	497 (98%)	9 (2%)	59 55		

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	58	LYS
1	A	74	PHE
1	A	172	ILE
1	A	219	PHE
1	A	239	ASN
1	В	39	LYS
1	В	74	PHE
1	В	193	LYS



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			- 0
Mol	Chain	Res	Type
1	В	219	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	18	GLN
1	A	80	GLN
1	A	94	ASN
1	A	140	HIS
1	A	239	ASN
1	В	67	ASN
1	В	80	GLN
1	В	94	ASN

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 monosaccharides in this entry.

5.6 Ligand geometry (i)

6 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with |Z| > 2 is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).



Mol	Tuno	Chain	Chain Res		Bo	ond leng	ths	В	ond ang	gles
MIOI	Type	Chain	nes	Link	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	MPO	В	303	-	13,13,13	1.97	1 (7%)	17,17,17	2.74	9 (52%)
2	1YJ	A	301	-	31,34,34	1.53	1 (3%)	39,47,47	1.96	18 (46%)
3	T3Q	A	302	-	34,37,37	1.37	5 (14%)	52,57,57	1.99	14 (26%)
2	1YJ	В	301	-	31,34,34	1.34	2 (6%)	39,47,47	2.03	13 (33%)
5	1YF	В	302	-	36,39,39	2.30	7 (19%)	52,59,59	1.81	11 (21%)
4	MPO	A	303	-	13,13,13	1.95	1 (7%)	17,17,17	2.60	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 Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	MPO	В	303	-	=	1/7/15/15	0/1/1/1
2	1YJ	A	301	-	ı	1/22/31/31	0/3/3/3
3	T3Q	A	302	_	-	2/21/53/53	0/3/3/3
2	1YJ	В	301	-	-	4/22/31/31	0/3/3/3
5	1YF	В	302	-	-	3/24/56/56	0/3/3/3
4	MPO	A	303	-	=	1/7/15/15	0/1/1/1

All (17) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$\operatorname{Observed}(\operatorname{\AA})$	Ideal(Å)
5	В	302	1YF	O2-C2	8.44	1.38	1.23
2	A	301	1YJ	O4-C4	6.83	1.36	1.23
5	В	302	1YF	O4-C4	6.78	1.36	1.23
4	В	303	MPO	C1-S1	-6.55	1.68	1.77
4	A	303	MPO	C1-S1	-6.51	1.68	1.77
2	В	301	1YJ	O4-C4	5.22	1.33	1.23
5	В	302	1YF	C1F-N3Q	4.45	1.48	1.33
3	A	302	T3Q	C6-N1	-3.79	1.31	1.38
5	В	302	1YF	C4-C5	-3.48	1.39	1.44
3	A	302	T3Q	C4-C5	-3.25	1.39	1.44
5	В	302	1YF	C6-C5	3.23	1.39	1.34
3	A	302	T3Q	O2-C2	2.68	1.27	1.23
3	A	302	T3Q	C6-C5	-2.67	1.30	1.34
5	В	302	1YF	O1F-C1F	2.45	1.30	1.22
2	В	301	1YJ	OX2-C	-2.25	1.23	1.30
3	A	302	T3Q	C2-N1	-2.03	1.35	1.38
5	В	302	1YF	C2-N1	-2.01	1.35	1.38



All (72) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^o)$	$\mathrm{Ideal}(^{o})$
4	В	303	MPO	C7-N1-C4	6.53	123.53	108.83
3	A	302	T3Q	C4-N3-C2	-6.12	119.43	127.35
5	В	302	1YF	C5-C4-N3	6.01	120.44	115.31
3	A	302	T3Q	O3A-PB-O3B	-5.77	90.85	102.48
4	A	303	MPO	C3-N1-C7	5.38	124.99	111.23
2	В	301	1YJ	O11-C11-C1'	-4.94	112.13	120.94
4	В	303	MPO	O1-S1-C1	4.67	112.53	106.92
4	A	303	MPO	C6-C7-N1	-4.56	103.19	110.10
4	A	303	MPO	C7-N1-C4	4.45	118.84	108.83
2	A	301	1YJ	C7-C6-C9	4.44	120.96	112.64
5	В	302	1YF	C4-N3-C2	-4.43	121.61	127.35
5	В	302	1YF	O3A-PB-O3B	-4.42	93.58	102.48
3	A	302	T3Q	O4Q-C4Q-C3Q	-4.29	102.52	110.22
2	В	301	1YJ	C1'-C11-N	4.19	125.10	117.06
4	В	303	MPO	O3-S1-C1	4.11	112.42	105.77
5	В	302	1YF	O4-C4-C5	-3.85	120.44	124.90
2	В	301	1YJ	C2-N1-C8A	3.68	120.08	113.43
3	A	302	T3Q	O4-C4-C5	-3.60	120.73	124.90
4	A	303	MPO	O1-S1-C1	3.59	111.24	106.92
3	A	302	T3Q	C5-C4-N3	3.58	118.37	115.31
3	A	302	T3Q	N3-C2-N1	3.54	119.59	114.89
5	В	302	1YF	C5-C6-N1	-3.31	119.93	123.34
2	A	301	1YJ	CB-CA-N	-3.28	104.25	110.88
2	В	301	1YJ	C4A-C4-N3	3.26	122.01	112.31
4	В	303	MPO	C3-N1-C7	3.20	119.42	111.23
3	A	302	T3Q	C6-C5-C4	3.17	120.68	118.03
2	A	301	1YJ	O11-C11-C1'	-3.13	115.35	120.94
4	В	303	MPO	C6-C7-N1	2.96	114.59	110.10
2	В	301	1YJ	CB-CA-C	-2.95	103.24	110.35
2	В	301	1YJ	O4-C4-C4A	-2.91	120.58	127.24
3	A	302	T3Q	O2-C2-N3	-2.90	116.10	121.50
3	A	302	T3Q	O1B-PB-O2B	2.85	126.32	112.24
2	A	301	1YJ	C3'-C4'-N10	-2.84	115.09	120.97
2	A	301	1YJ	C-CA-N	2.82	117.22	110.55
2	A	301	1YJ	OE1-CD-CG	-2.82	114.04	123.08
3	A	302	T3Q	C5-C6-N1	-2.81	120.44	123.34
2	В	301	1YJ	OE2-CD-CG	2.80	123.03	114.03
2	В	301	1YJ	C2-N3-C4	-2.80	119.99	125.10
2	A	301	1YJ	C1'-C11-N	2.77	122.37	117.06
5	В	302	1YF	O5Q-C1Q-C2Q	2.73	116.13	110.35
3	A	302	T3Q	O2A-PA-O1A	2.69	125.52	112.24
2	A	301	1YJ	C5'-C4'-N10	2.67	126.50	120.97



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Mol	Chain	Res	Type	Atoms	Z	$\operatorname{Observed}({}^o)$	$\operatorname{Ideal}({}^{o})$
2	В	301	1YJ	C9-C6-N5	-2.66	103.11	108.83
2	A	301	1YJ	C6-C7-N8	2.64	115.63	110.68
5	В	302	1YF	N3-C2-N1	2.61	118.35	114.89
2	A	301	1YJ	CB-CA-C	-2.47	104.40	110.35
4	В	303	MPO	C5-C4-N1	-2.46	106.37	110.10
2	В	301	1YJ	C-CA-N	2.45	116.34	110.55
2	A	301	1YJ	C2-N3-C4	-2.42	120.69	125.10
2	A	301	1YJ	C2-N1-C8A	2.40	117.77	113.43
4	В	303	MPO	O2-S1-O1	-2.40	105.65	113.95
3	A	302	T3Q	O4Q-C4Q-C5Q	2.36	114.90	109.67
2	A	301	1YJ	OX2-C-CA	2.29	121.02	113.40
4	A	303	MPO	O3-S1-C1	2.28	109.46	105.77
4	В	303	MPO	C3-N1-C4	2.26	117.01	111.23
2	A	301	1YJ	OE2-CD-CG	2.25	121.27	114.03
2	В	301	1YJ	C6-C7-N8	2.25	114.89	110.68
2	A	301	1YJ	C9-C6-N5	-2.23	104.03	108.83
2	A	301	1YJ	O4-C4-N3	-2.18	115.93	120.12
2	В	301	1YJ	CB-CA-N	-2.18	106.47	110.88
3	A	302	T3Q	C5Q-C4Q-C3Q	-2.17	107.22	110.51
5	В	302	1YF	O1B-PB-O2B	2.14	122.82	112.24
2	A	301	1YJ	C4A-C4-N3	2.12	118.61	112.31
4	A	303	MPO	O2-S1-O1	-2.11	106.65	113.95
4	A	303	MPO	C2-C1-S1	2.09	116.45	113.25
5	В	302	1YF	O1B-PB-O3B	-2.08	98.56	106.78
2	В	301	1YJ	CG-CB-CA	2.07	117.02	113.16
5	В	302	1YF	O5Q-C1Q-O3B	-2.06	108.67	111.36
3	A	302	T3Q	C6Q-C5Q-C4Q	2.06	116.87	113.07
4	В	303	MPO	O4-C5-C4	-2.06	107.27	111.80
2	A	301	1YJ	OX1-C-CA	-2.05	115.53	122.26
5	В	302	1YF	O4Q-C4Q-C3Q	-2.02	105.57	109.66

There are no chirality outliers.

All (12) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	301	1YJ	C7-C6-C9-N10
3	A	302	T3Q	PB-O3A-PA-O5'
5	В	302	1YF	O1F-C1F-N3Q-C3Q
4	A	303	MPO	C2-C3-N1-C7
4	В	303	MPO	C2-C3-N1-C4
5	В	302	1YF	PB-O3A-PA-O5'
3	A	302	T3Q	C1Q-O3B-PB-O3A



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Mol	Chain	Res	Type	Atoms
2	В	301	1YJ	C3'-C4'-N10-C9
5	В	302	1YF	C1Q-O3B-PB-O3A
2	В	301	1YJ	OE1-CD-CG-CB
2	В	301	1YJ	C-CA-CB-CG
2	В	301	1YJ	OE2-CD-CG-CB

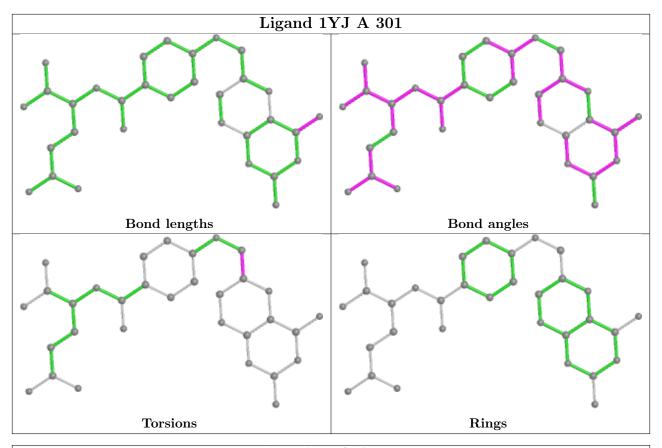
There are no ring outliers.

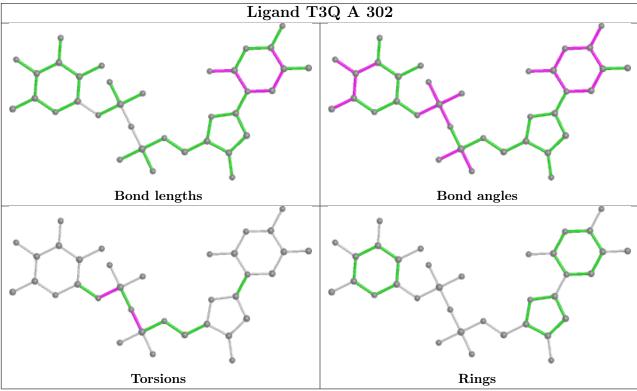
2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	В	303	MPO	1	0
5	В	302	1YF	1	0

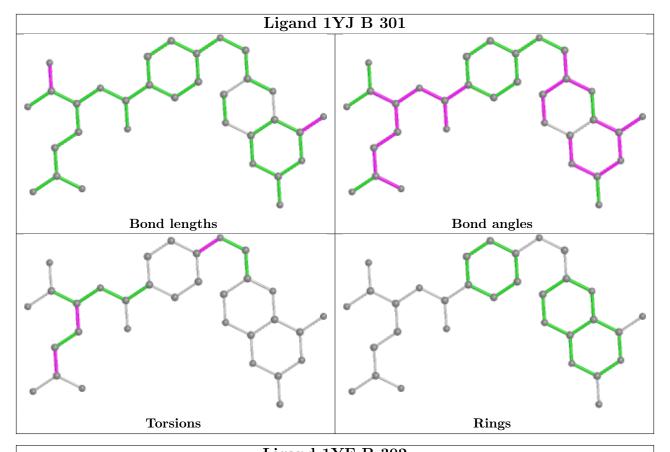
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less then 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.

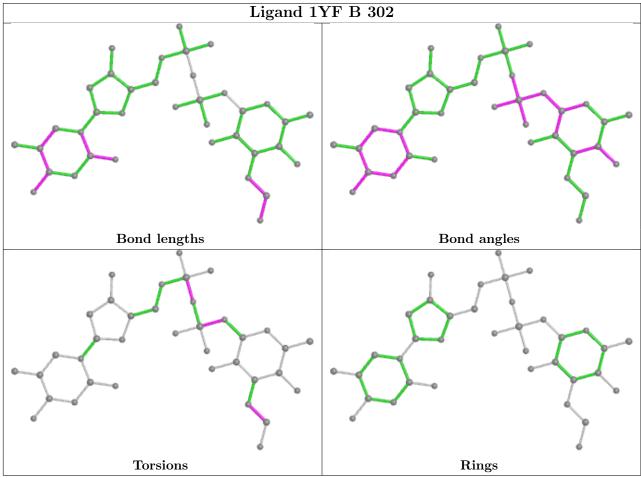














5.7 Other polymers (i)

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, 95^{th} 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	$\langle { m RSRZ} \rangle$	$\#\mathrm{RSRZ}{>}2$			$OWAB(A^2)$	Q < 0.9
1	A	271/274 (98%)	-0.61	1 (0%)	92	93	8, 16, 30, 55	0
1	В	271/274 (98%)	-0.55	1 (0%)	92	93	8, 15, 33, 63	0
All	All	542/548 (98%)	-0.58	2 (0%)	92	93	8, 16, 31, 63	0

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	270	ASN	3.0
1	В	-1	GLY	2.1

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 monosaccharides 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, 95^{th} 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.

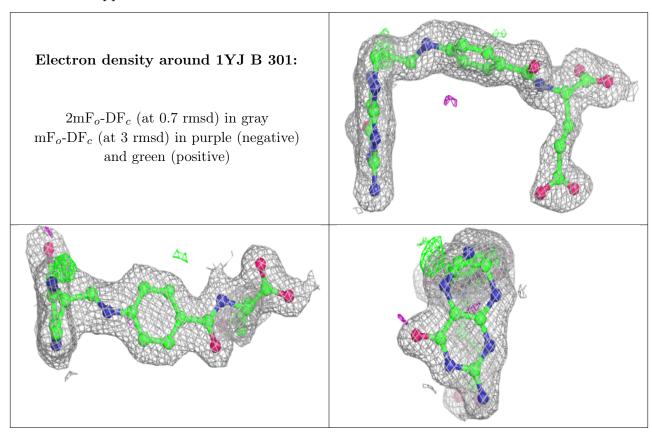
Mol	Type	Chain	Res	Atoms	RSCC	RSR	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{ ilde{A}}^2)$	Q<0.9
4	MPO	A	303	13/13	0.88	0.16	40,45,63,70	0
4	MPO	В	303	13/13	0.92	0.17	37,46,53,56	0
2	1YJ	В	301	32/32	0.94	0.10	14,21,34,44	0



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Mol	Type	Chain	Res	Atoms	RSCC	RSR	$\mathbf{B} ext{-}\mathbf{factors}(\mathring{\mathbf{A}}^2)$	Q < 0.9
2	1YJ	A	301	32/32	0.95	0.09	12,19,33,39	0
3	T3Q	A	302	35/35	0.97	0.09	14,17,24,28	0
5	1YF	В	302	37/37	0.97	0.08	12,16,25,30	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.

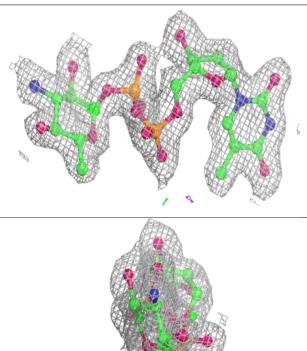


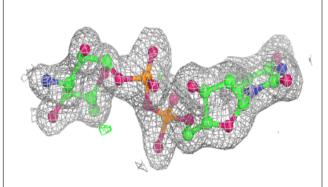


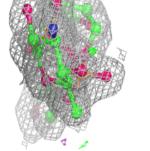
Electron density around 1YJ A 301: $2 \mathrm{mF}_o\text{-}\mathrm{DF}_c$ (at 0.7 rmsd) in gray ${\rm mF}_o\text{-}{\rm DF}_c$ (at 3 rmsd) in purple (negative) and green (positive)

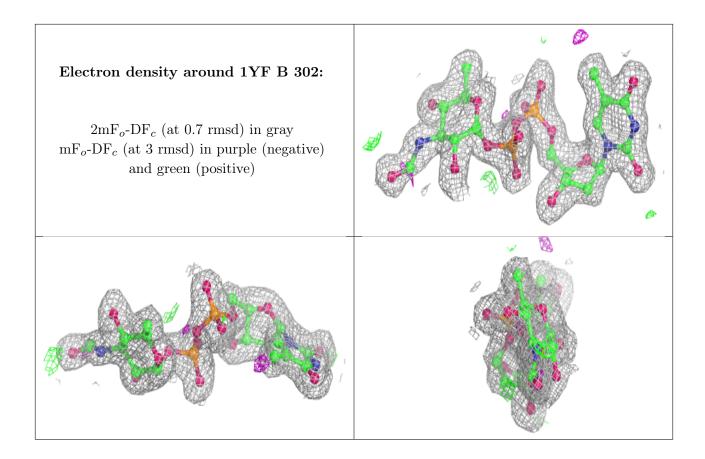
Electron density around T3Q A 302:

 $2 \mathrm{mF}_o\text{-}\mathrm{DF}_c$ (at 0.7 rmsd) in gray $\mathrm{mF}_o\text{-}\mathrm{DF}_c$ (at 3 rmsd) in purple (negative) and green (positive)









6.5 Other polymers (i)

There are no such residues in this entry.

