



Full wwPDB X-ray Structure Validation Report ⓘ

May 18, 2020 – 02:13 am BST

PDB ID : 1NZC
Title : The high resolution structures of RmlC from *Streptococcus suis* in complex with dTDP-D-xylose
Authors : Dong, C.; Major, L.L.; Allen, A.; Blankenfeldt, W.; Maskell, D.; Naismith, J.H.
Deposited on : 2003-02-17
Resolution : 1.80 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

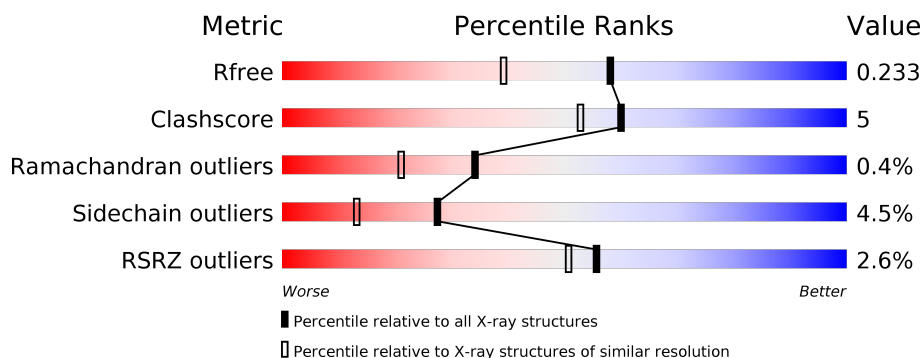
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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	5950 (1.80-1.80)
Clashscore	141614	6793 (1.80-1.80)
Ramachandran outliers	138981	6697 (1.80-1.80)
Sidechain outliers	138945	6696 (1.80-1.80)
RSRZ outliers	127900	5850 (1.80-1.80)

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	197	<div> <div>%</div> <div> <div></div> <div>86%</div> <div>11%</div> <div>...</div> </div> </div>
1	B	197	<div> <div>7%</div> <div> <div></div> <div>82%</div> <div>15%</div> <div>...</div> </div> </div>
1	C	197	<div> <div>%</div> <div> <div></div> <div>86%</div> <div>11%</div> <div>...</div> </div> </div>
1	D	197	<div> <div>%</div> <div> <div></div> <div>86%</div> <div>11%</div> <div>..</div> </div> </div>

2 Entry composition [i](#)

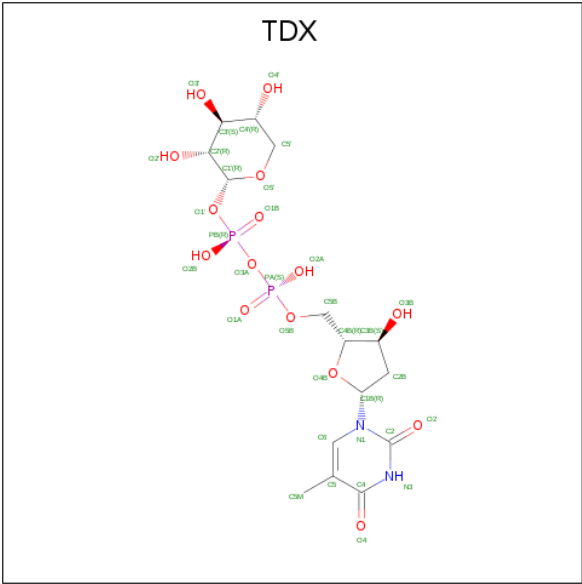
There are 4 unique types of molecules in this entry. The entry contains 7316 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 dTDP-6-deoxy-D-xylo-4-hexulose 3,5-epimerase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	196	Total 1584	1026	259	297	2	0	0	0
1	B	196	Total 1584	1026	259	297	2	0	0	0
1	C	194	Total 1568	1017	257	292	2	0	0	0
1	D	194	Total 1568	1017	257	292	2	0	0	0

- Molecule 2 is THYMIDINE-5'-DIPHOSPHO-BETA-D-XYLOSE (three-letter code: TDX) (formula: C₁₅H₂₄N₂O₁₅P₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
2	A	1	Total 34	15	2	15	2	0	0
2	B	1	Total 34	15	2	15	2	1	0

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	C	1	Total	C	N	O	P	0	0
			34	15	2	15	2		
2	D	1	Total	C	N	O	P	0	0
			34	15	2	15	2		

- Molecule 3 is NICKEL (II) ION (three-letter code: NI) (formula: Ni).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Ni	0	0
			1	1		
3	C	1	Total	Ni	0	0
			1	1		

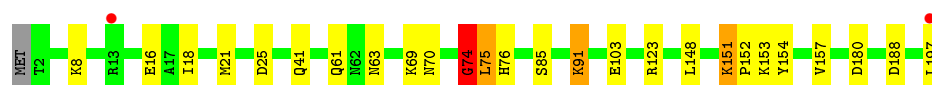
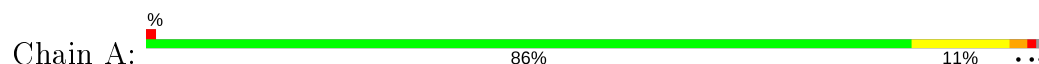
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	225	Total	O	0	0
			225	225		
4	B	198	Total	O	0	0
			198	198		
4	C	246	Total	O	0	0
			246	246		
4	D	205	Total	O	0	0
			205	205		

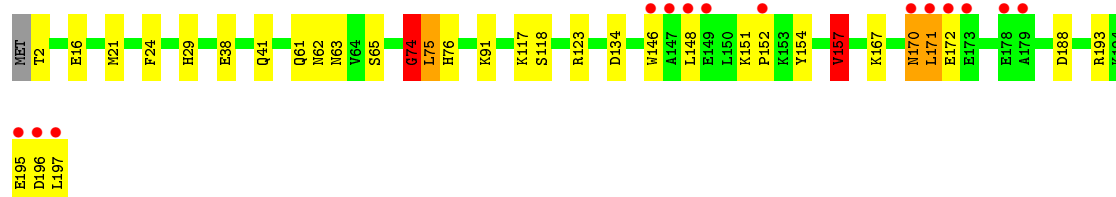
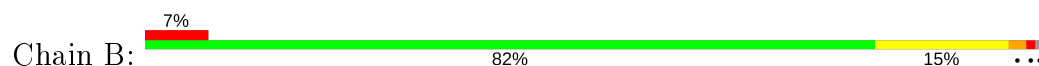
3 Residue-property plots [i](#)

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.

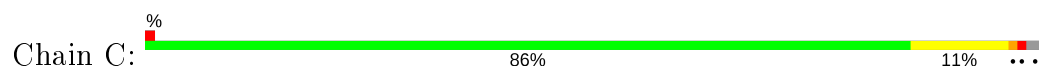
- Molecule 1: dTDP-6-deoxy-D-xylo-4-hexulose 3,5-epimerase



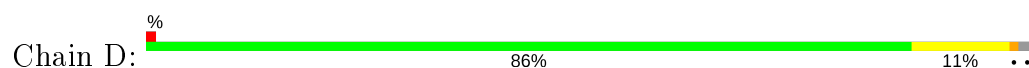
- Molecule 1: dTDP-6-deoxy-D-xylo-4-hexulose 3,5-epimerase



- Molecule 1: dTDP-6-deoxy-D-xylo-4-hexulose 3,5-epimerase



- Molecule 1: dTDP-6-deoxy-D-xylo-4-hexulose 3,5-epimerase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	50.85Å 140.87Å 53.70Å 90.00° 92.72° 90.00°	Depositor
Resolution (Å)	70.71 – 1.80 50.79 – 1.80	Depositor EDS
% Data completeness (in resolution range)	91.8 (70.71-1.80) 91.8 (50.79-1.80)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	0.12	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.44 (at 1.79Å)	Xtriage
Refinement program	REFMAC 5.1.19	Depositor
R, R_{free}	0.164 , 0.216 0.193 , 0.233	Depositor DCC
R_{free} test set	3217 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	20.0	Xtriage
Anisotropy	0.114	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 46.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.032 for h,-k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7316	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.30% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: NI, TDX

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	0.61	0/1629	0.82	5/2209 (0.2%)
1	B	0.67	2/1629 (0.1%)	0.87	4/2209 (0.2%)
1	C	0.63	0/1613	0.82	5/2187 (0.2%)
1	D	0.59	0/1613	0.85	9/2187 (0.4%)
All	All	0.63	2/6484 (0.0%)	0.84	23/8792 (0.3%)

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	A	0	2
1	B	0	1
1	C	0	1
All	All	0	4

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	62	ASN	CG-ND2	-5.74	1.18	1.32
1	B	62	ASN	CG-OD1	-5.16	1.12	1.24

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	74	GLY	CA-C-N	-7.54	100.60	117.20
1	D	74	GLY	CA-C-N	-7.32	101.10	117.20
1	B	188	ASP	CB-CG-OD2	7.26	124.83	118.30
1	D	74	GLY	O-C-N	7.15	134.14	122.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	74	GLY	CA-C-N	-6.85	102.13	117.20
1	D	74	GLY	C-N-CA	6.82	138.76	121.70
1	C	74	GLY	O-C-N	6.50	133.10	122.70
1	A	74	GLY	C-N-CA	6.40	137.70	121.70
1	D	188	ASP	CB-CG-OD2	6.28	123.95	118.30
1	C	74	GLY	C-N-CA	6.27	137.38	121.70
1	D	75	LEU	CB-CA-C	6.24	122.06	110.20
1	A	74	GLY	O-C-N	6.18	132.58	122.70
1	A	188	ASP	CB-CG-OD2	5.94	123.64	118.30
1	C	75	LEU	CB-CA-C	5.93	121.47	110.20
1	D	196	ASP	CB-CG-OD2	5.56	123.31	118.30
1	D	180	ASP	CB-CG-OD2	5.52	123.27	118.30
1	A	180	ASP	CB-CG-OD2	5.49	123.24	118.30
1	D	81	ASP	CB-CG-OD2	5.44	123.20	118.30
1	B	134	ASP	CB-CG-OD2	5.41	123.17	118.30
1	D	74	GLY	N-CA-C	5.33	126.41	113.10
1	B	157	VAL	CB-CA-C	-5.27	101.39	111.40
1	C	180	ASP	CB-CG-OD2	5.13	122.92	118.30
1	B	74	GLY	CA-C-N	-5.11	105.96	117.20

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	74	GLY	Mainchain,Peptide
1	B	74	GLY	Peptide
1	C	74	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	1584	0	1525	17	1
1	B	1584	0	1525	17	2
1	C	1568	0	1512	16	0
1	D	1568	0	1512	10	0
2	A	34	0	22	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	34	0	22	0	0
2	C	34	0	22	1	0
2	D	34	0	22	0	0
3	B	1	0	0	1	0
3	C	1	0	0	0	0
4	A	225	0	0	8	6
4	B	198	0	0	6	2
4	C	246	0	0	3	4
4	D	205	0	0	0	1
All	All	7316	0	6162	58	8

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 (58) 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:C:61:GLN:HE21	1:C:63:ASN:HD21	1.08	0.99
1:D:61:GLN:HE21	1:D:63:ASN:HD21	1.15	0.94
1:B:61:GLN:HE21	1:B:63:ASN:HD21	1.07	0.93
1:B:196:ASP:OD2	4:B:9634:HOH:O	1.87	0.92
1:A:61:GLN:HE21	1:A:63:ASN:HD21	1.18	0.89
1:A:85:SER:OG	4:A:9077:HOH:O	1.58	0.82
1:B:76:HIS:ND1	4:B:9636:HOH:O	2.04	0.80
3:B:9603:NI:NI	4:B:9656:HOH:O	1.28	0.75
1:D:74:GLY:HA2	1:D:76:HIS:CE1	2.23	0.74
1:C:197:LEU:HD23	4:C:9813:HOH:O	1.89	0.71
1:D:46:LEU:HD23	1:D:50:PHE:O	1.92	0.69
1:B:170:ASN:O	1:B:172:GLU:N	2.27	0.67
1:C:74:GLY:HA2	1:C:76:HIS:CE1	2.29	0.67
1:A:197:LEU:HA	4:A:9121:HOH:O	1.95	0.66
1:A:123:ARG:HD3	4:A:9074:HOH:O	1.95	0.66
1:A:74:GLY:HA2	1:A:76:HIS:CE1	2.31	0.65
1:A:91:LYS:HG2	4:A:9003:HOH:O	1.99	0.63
1:B:151:LYS:HA	1:B:154:TYR:CD1	2.34	0.63
1:B:195:GLU:HG3	4:B:9654:HOH:O	2.00	0.61
1:C:75:LEU:HB2	1:C:157:VAL:O	2.00	0.61
1:C:61:GLN:NE2	1:C:63:ASN:HD21	1.90	0.60
1:C:197:LEU:C	4:C:9816:HOH:O	2.42	0.57
1:C:29:HIS:HD2	4:C:9754:HOH:O	1.87	0.57
1:A:153:LYS:HG2	4:A:9206:HOH:O	2.04	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:74:GLY:CA	1:D:76:HIS:CE1	2.88	0.57
1:A:8:LYS:HD3	1:A:25:ASP:HB3	1.89	0.55
1:C:76:HIS:ND1	2:C:9003:TDX:H3'	2.23	0.54
1:C:175:GLU:OE2	1:D:32:ASN:ND2	2.41	0.54
4:A:9072:HOH:O	1:B:29:HIS:HE1	1.91	0.54
1:D:159:TYR:OH	1:D:181:GLU:OE2	2.17	0.53
1:B:65:SER:HB2	4:B:9628:HOH:O	2.09	0.52
1:B:151:LYS:N	1:B:152:PRO:CD	2.73	0.52
1:D:61:GLN:NE2	1:D:63:ASN:HD21	1.97	0.51
1:C:123:ARG:HH11	1:C:197:LEU:HD22	1.77	0.49
1:B:61:GLN:NE2	1:B:63:ASN:HD21	1.91	0.48
1:B:151:LYS:HA	1:B:154:TYR:CE1	2.49	0.47
1:B:196:ASP:O	1:B:197:LEU:HB2	2.15	0.47
1:A:18:ILE:HB	1:A:21:MET:HE3	1.96	0.47
1:A:61:GLN:NE2	1:A:63:ASN:HD21	2.00	0.47
1:B:75:LEU:HB2	1:B:157:VAL:O	2.15	0.46
1:A:151:LYS:HA	1:A:154:TYR:CD2	2.51	0.46
1:A:63:ASN:HD22	1:B:38:GLU:HA	1.81	0.45
1:C:151:LYS:N	1:C:152:PRO:CD	2.79	0.45
1:D:176:VAL:CG2	1:D:181:GLU:HG2	2.47	0.45
1:D:75:LEU:HB2	1:D:157:VAL:O	2.17	0.44
1:A:16:GLU:HG2	4:A:9082:HOH:O	2.17	0.44
1:B:117:LYS:HE3	4:B:9787:HOH:O	2.18	0.43
1:A:151:LYS:N	1:A:152:PRO:CD	2.81	0.43
1:B:24:PHE:HB2	1:B:118:SER:OG	2.18	0.43
1:A:75:LEU:HB2	1:A:157:VAL:O	2.19	0.43
1:C:46:LEU:HB2	1:C:47:PRO:HD3	2.01	0.42
1:C:38:GLU:HA	1:D:63:ASN:HD22	1.84	0.42
1:A:8:LYS:NZ	4:A:9141:HOH:O	2.42	0.42
1:B:74:GLY:HA2	1:B:76:HIS:CE1	2.55	0.41
1:C:180:ASP:HA	1:C:183:HIS:CD2	2.56	0.41
1:C:24:PHE:HB2	1:C:118:SER:OG	2.21	0.41
1:A:69:LYS:O	1:A:70:ASN:HB2	2.21	0.40
1:C:176:VAL:CG1	1:C:180:ASP:HB2	2.52	0.40

All (8) 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)
1:B:2:THR:OG1	4:A:9019:HOH:O[1_655]	1.30	0.90

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:16:GLU:OE1	4:B:9697:HOH:O[1_455]	1.74	0.46
4:B:9654:HOH:O	4:D:9102:HOH:O[1_545]	1.79	0.41
4:A:9035:HOH:O	4:C:9823:HOH:O[2_546]	2.03	0.17
1:B:2:THR:N	4:A:9212:HOH:O[1_655]	2.06	0.14
4:A:9035:HOH:O	4:C:9689:HOH:O[2_546]	2.13	0.07
4:A:9035:HOH:O	4:C:9717:HOH:O[2_546]	2.17	0.03
4:A:9185:HOH:O	4:C:9715:HOH:O[2_546]	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	194/197 (98%)	189 (97%)	4 (2%)	1 (0%)	29	15
1	B	194/197 (98%)	190 (98%)	2 (1%)	2 (1%)	15	5
1	C	192/197 (98%)	185 (96%)	7 (4%)	0	100	100
1	D	192/197 (98%)	183 (95%)	9 (5%)	0	100	100
All	All	772/788 (98%)	747 (97%)	22 (3%)	3 (0%)	34	21

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	171	LEU
1	A	75	LEU
1	B	75	LEU

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	168/169 (99%)	163 (97%)	5 (3%)	41	27
1	B	168/169 (99%)	156 (93%)	12 (7%)	14	5
1	C	166/169 (98%)	161 (97%)	5 (3%)	41	27
1	D	166/169 (98%)	158 (95%)	8 (5%)	25	11
All	All	668/676 (99%)	638 (96%)	30 (4%)	27	13

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	41	GLN
1	A	91	LYS
1	A	103	GLU
1	A	148	LEU
1	A	151	LYS
1	B	16	GLU
1	B	21	MET
1	B	41	GLN
1	B	91	LYS
1	B	123	ARG
1	B	146	TRP
1	B	148	LEU
1	B	157	VAL
1	B	167	LYS
1	B	170	ASN
1	B	171	LEU
1	B	193	ARG
1	C	41	GLN
1	C	91	LYS
1	C	148	LEU
1	C	167	LYS
1	C	196	ASP
1	D	16	GLU
1	D	41	GLN
1	D	91	LYS
1	D	148	LEU
1	D	149	GLU
1	D	150	LEU
1	D	163	SER
1	D	197	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	39	ASN
1	A	63	ASN
1	B	29	HIS
1	B	62	ASN
1	B	63	ASN
1	B	170	ASN
1	C	29	HIS
1	C	63	ASN
1	D	29	HIS
1	D	63	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 6 ligands modelled in this entry, 2 are 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	TDX	B	9001	-	30,36,36	0.91	1 (3%)	40,55,55	1.97	5 (12%)
2	TDX	D	9004	-	30,36,36	0.90	1 (3%)	40,55,55	1.88	3 (7%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	TDX	A	9002	-	30,36,36	0.91	1 (3%)	40,55,55	1.60	3 (7%)
2	TDX	C	9003	-	30,36,36	0.78	1 (3%)	40,55,55	1.84	2 (5%)

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	TDX	B	9001	-	-	4/18/50/50	0/3/3/3
2	TDX	D	9004	-	-	3/18/50/50	0/3/3/3
2	TDX	A	9002	-	-	3/18/50/50	0/3/3/3
2	TDX	C	9003	-	-	3/18/50/50	0/3/3/3

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	9001	TDX	C4-C5	3.62	1.49	1.41
2	A	9002	TDX	C4-C5	3.58	1.49	1.41
2	D	9004	TDX	C4-C5	3.17	1.48	1.41
2	C	9003	TDX	C4-C5	2.55	1.46	1.41

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	9004	TDX	C4-N3-C2	10.02	123.60	115.14
2	C	9003	TDX	C4-N3-C2	9.57	123.22	115.14
2	B	9001	TDX	C4-N3-C2	8.51	122.33	115.14
2	A	9002	TDX	C4-N3-C2	8.11	121.99	115.14
2	B	9001	TDX	O3'-C3'-C4'	5.71	120.93	109.99
2	B	9001	TDX	O3'-C3'-C2'	-2.92	103.60	110.35
2	D	9004	TDX	C5'-C4'-C3'	2.63	112.90	109.67
2	D	9004	TDX	C2B-C1B-N1	2.59	120.23	114.27
2	C	9003	TDX	O2B-PB-O1B	2.42	124.19	112.24
2	B	9001	TDX	O5'-C1'-O1'	-2.23	108.45	111.36
2	A	9002	TDX	O2B-PB-O1B	2.10	122.65	112.24
2	B	9001	TDX	C5-C6-N1	-2.03	120.00	122.19
2	A	9002	TDX	C4'-C3'-C2'	2.02	114.39	110.89

There are no chirality outliers.

All (13) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	9001	TDX	O4B-C1B-N1-C6
2	B	9001	TDX	C5B-O5B-PA-O1A
2	B	9001	TDX	C5B-O5B-PA-O2A
2	A	9002	TDX	O4B-C1B-N1-C6
2	D	9004	TDX	O4B-C1B-N1-C6
2	C	9003	TDX	O4B-C1B-N1-C6
2	D	9004	TDX	PA-O3A-PB-O1B
2	A	9002	TDX	PA-O3A-PB-O1B
2	C	9003	TDX	PA-O3A-PB-O1B
2	B	9001	TDX	C5B-O5B-PA-O3A
2	A	9002	TDX	PA-O3A-PB-O2B
2	D	9004	TDX	PA-O3A-PB-O2B
2	C	9003	TDX	PA-O3A-PB-O2B

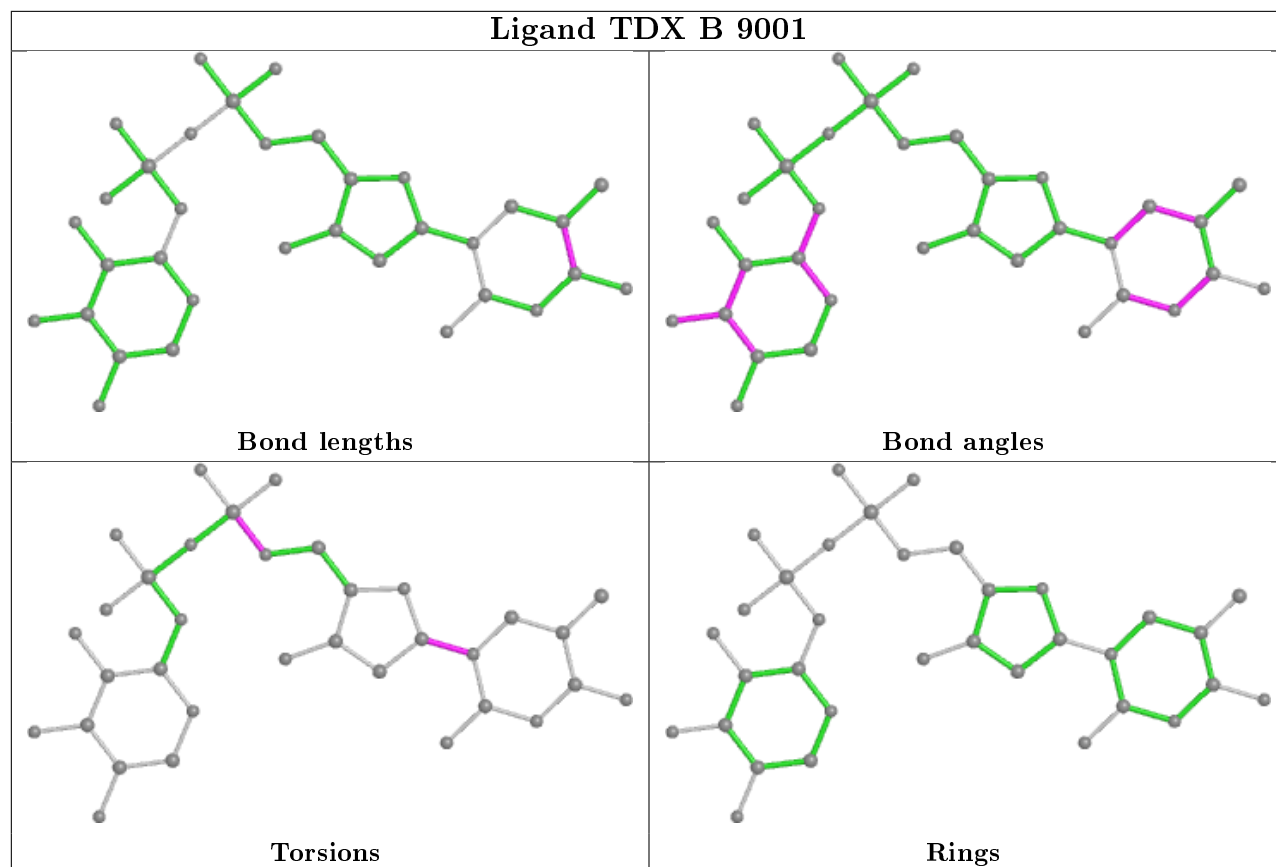
There are no ring outliers.

1 monomer is involved in 1 short contact:

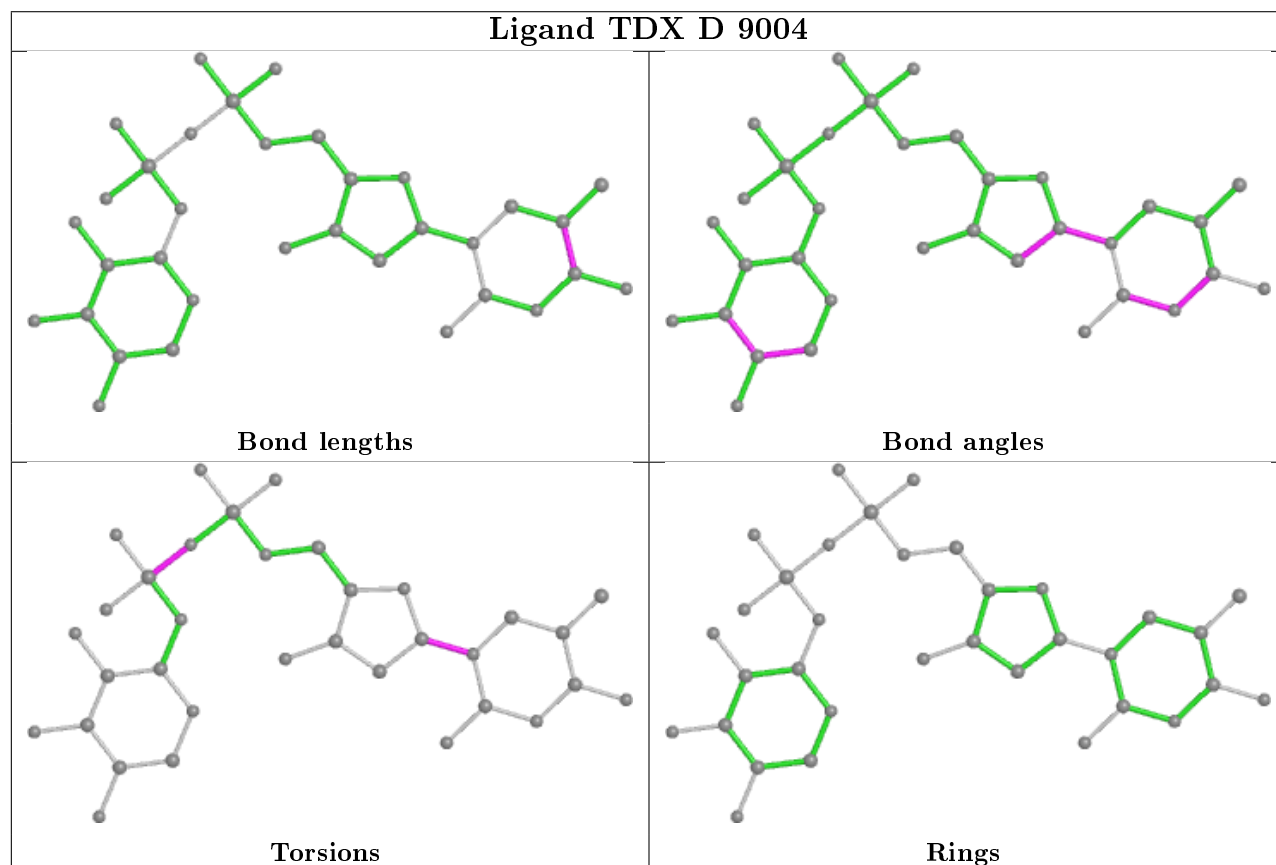
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	9003	TDX	1	0

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.

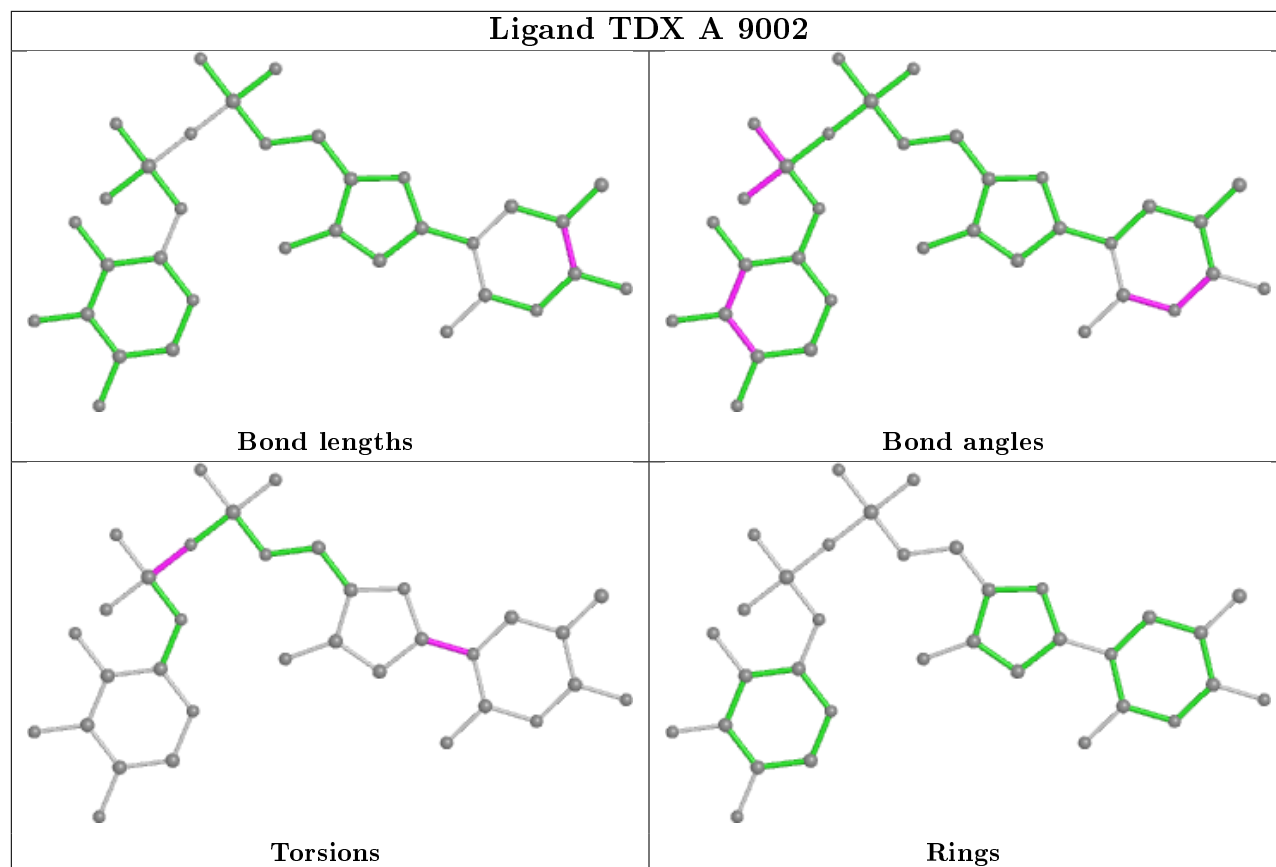
Ligand TDX B 9001



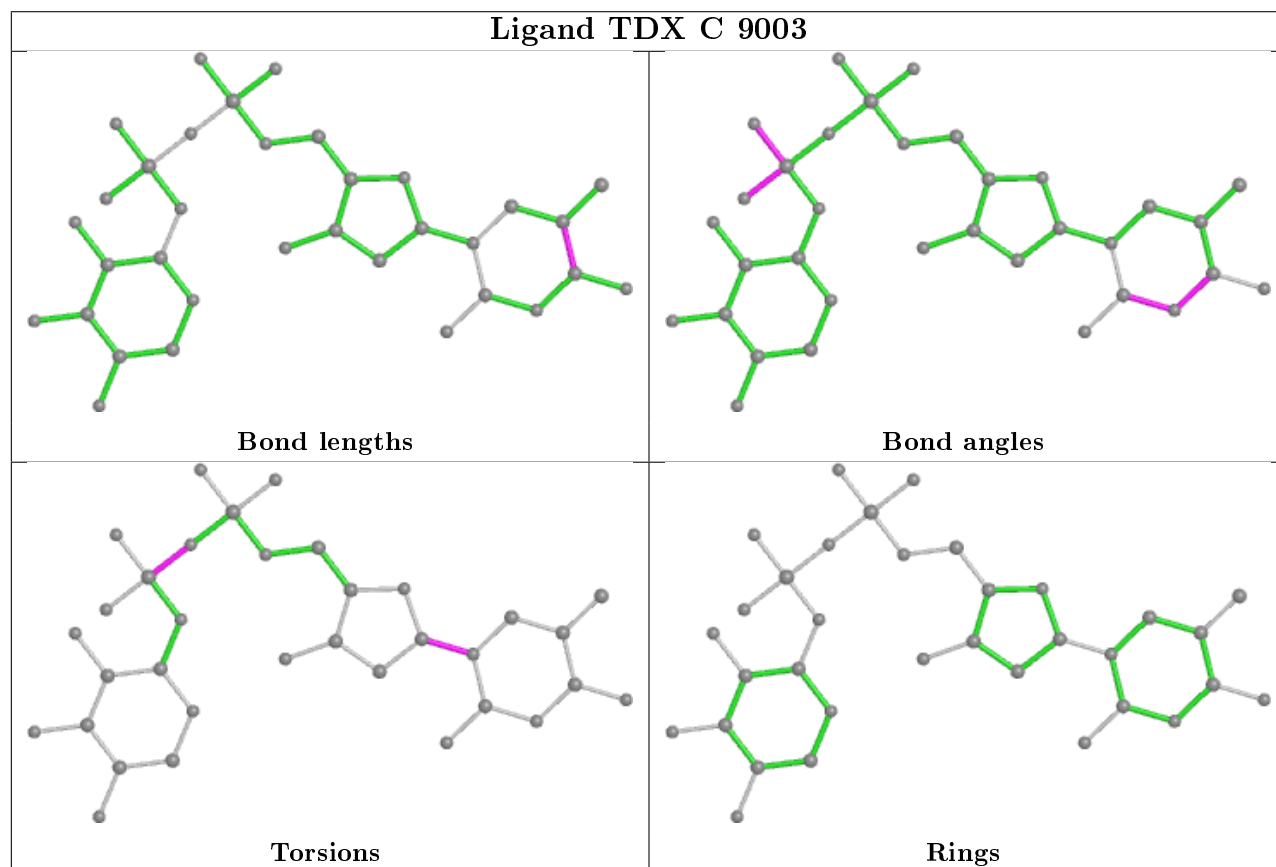
Ligand TDX D 9004



Ligand TDX A 9002



Ligand TDX C 9003



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	A	196/197 (99%)	-0.14	2 (1%) 82 80	12, 17, 25, 38	0
1	B	196/197 (99%)	0.25	14 (7%) 16 12	15, 18, 27, 38	0
1	C	194/197 (98%)	-0.11	2 (1%) 82 80	14, 18, 26, 38	0
1	D	194/197 (98%)	-0.15	2 (1%) 82 80	15, 18, 27, 37	0
All	All	780/788 (98%)	-0.04	20 (2%) 56 51	12, 18, 27, 38	0

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	148	LEU	6.0
1	B	146	TRP	3.7
1	B	172	GLU	3.6
1	B	197	LEU	3.6
1	C	197	LEU	3.5
1	B	152	PRO	3.3
1	B	171	LEU	3.0
1	B	149	GLU	3.0
1	D	195	GLU	2.7
1	B	196	ASP	2.6
1	B	179	ALA	2.6
1	B	173	GLU	2.5
1	B	147	ALA	2.4
1	A	197	LEU	2.3
1	B	170	ASN	2.3
1	A	13	ARG	2.3
1	B	178	GLU	2.2
1	B	195	GLU	2.2
1	D	16	GLU	2.1
1	C	196	ASP	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](#)

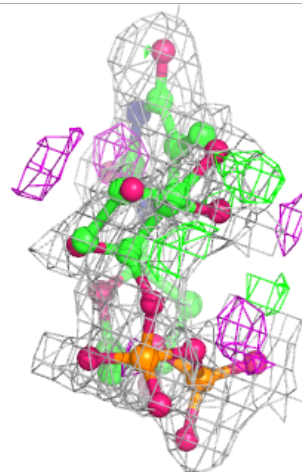
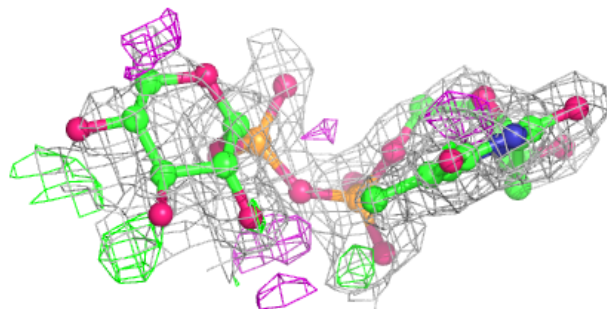
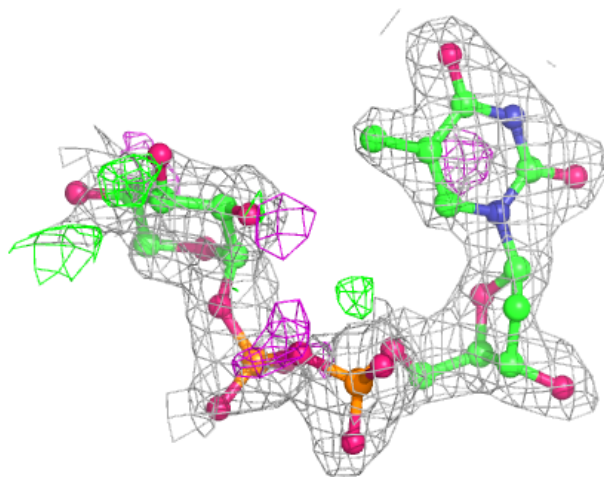
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	TDX	B	9001	34/34	0.77	0.21	41,64,67,68	1
3	NI	B	9603	1/1	0.89	0.34	128,128,128,128	0
2	TDX	A	9002	34/34	0.96	0.11	20,22,24,25	0
2	TDX	D	9004	34/34	0.96	0.10	20,23,26,27	0
2	TDX	C	9003	34/34	0.96	0.10	19,23,27,27	0
3	NI	C	9602	1/1	0.99	0.11	49,49,49,49	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.

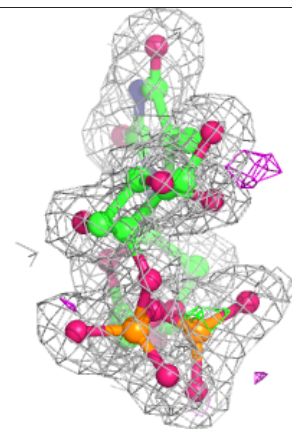
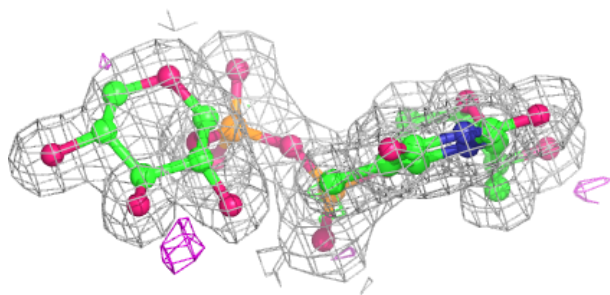
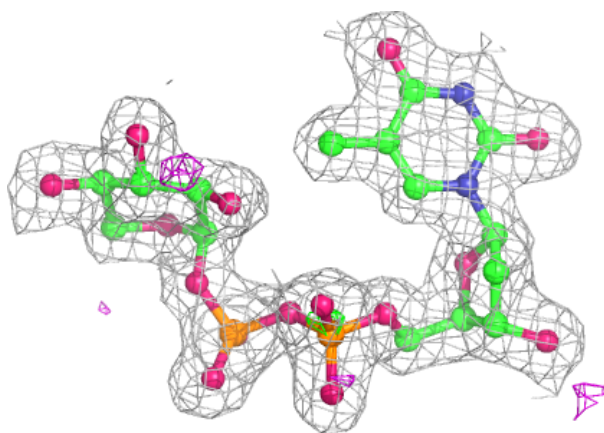
Electron density around TDX B 9001:

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



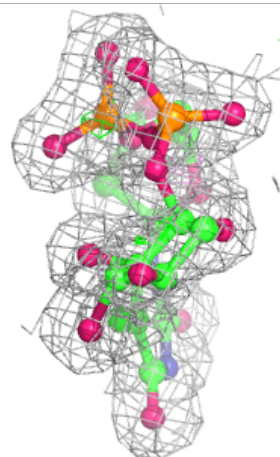
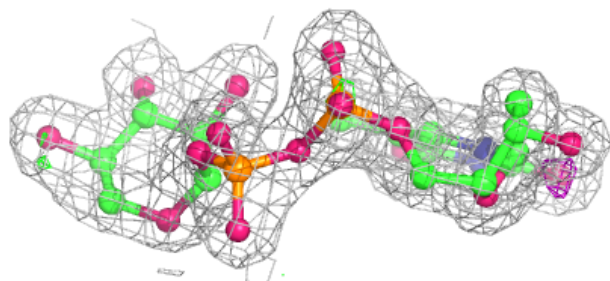
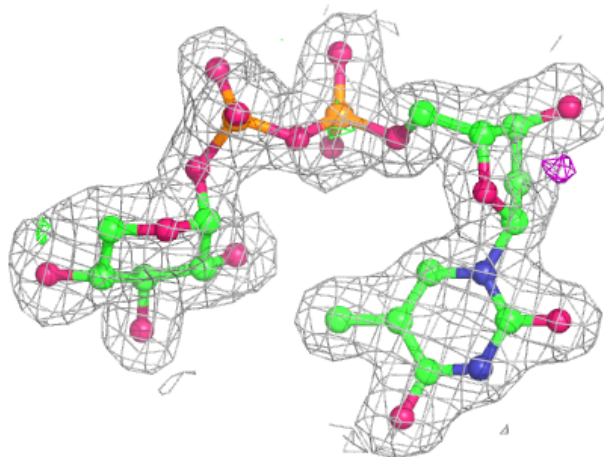
Electron density around TDX A 9002:

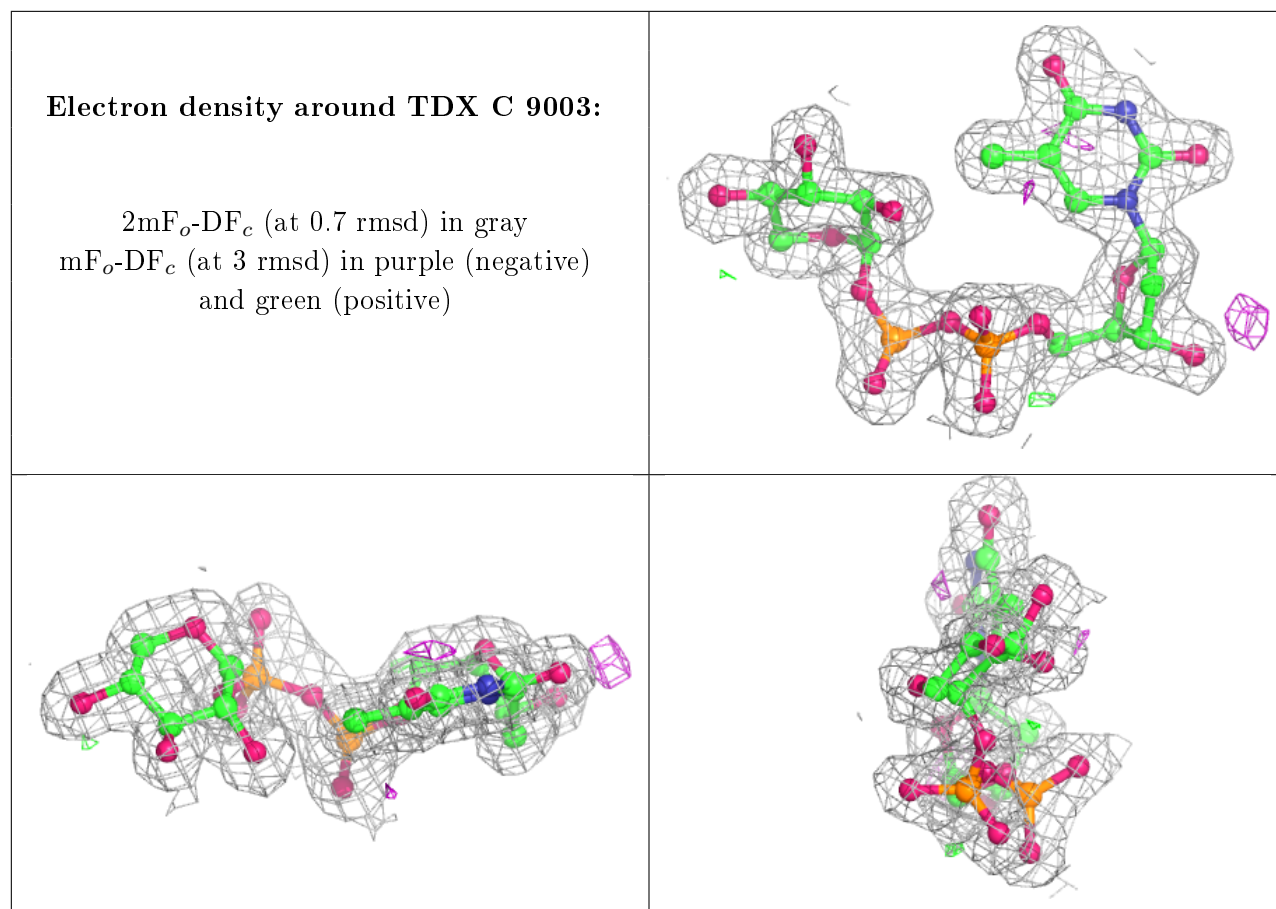
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around TDX D 9004:

$2mF_o-DF_c$ (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.