



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 22, 2026 – 03:11 am BST

PDB ID : 2BVL / pdb_00002bvl
Title : Crystal structure of the catalytic domain of toxin B from *Clostridium difficile* in complex with UDP, Glc and manganese ion
Authors : Reinert, D.J.; Jank, T.; Aktories, K.; Schulz, G.E.
Deposited on : 2005-06-30
Resolution : 2.20 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.015 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.50

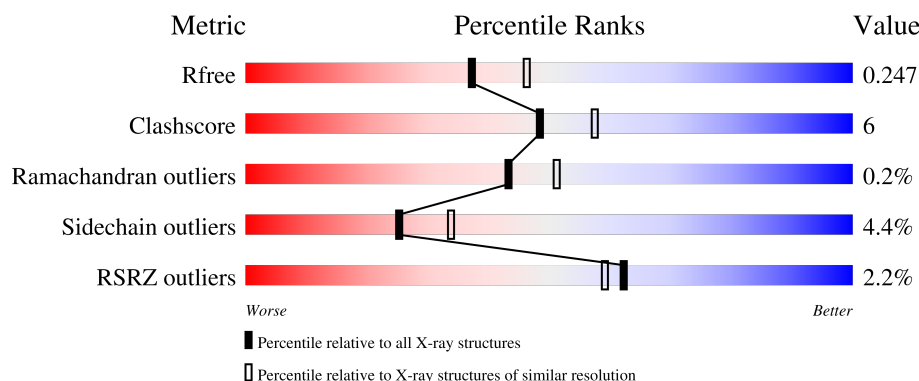
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

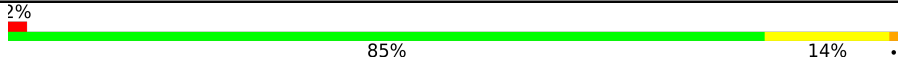
The reported resolution of this entry is 2.20 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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

Mol	Chain	Length	Quality of chain
1	A	543	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
6	TBR	A	1549	-	-	X	-

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 4902 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 TOXIN B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	543	Total	C	N	O	S	0	0	0
			4433	2810	720	885	18			

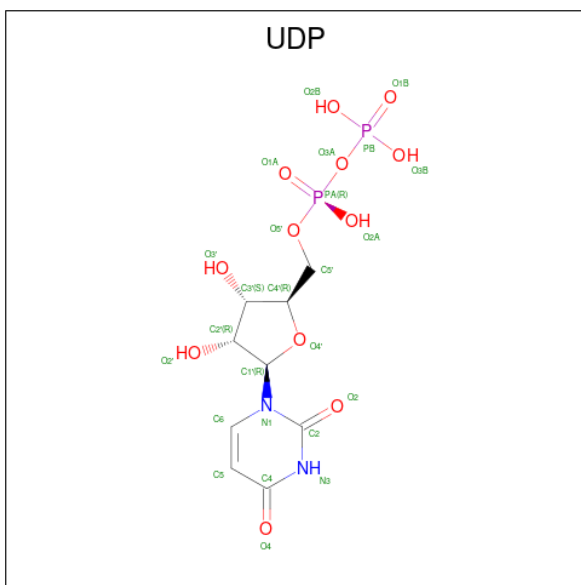
There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	59	CYS	TYR	engineered mutation	UNP P18177
A	244	GLY	GLU	engineered mutation	UNP P18177
A	542	ALA	SER	engineered mutation	UNP P18177

- Molecule 2 is MANGANESE (II) ION (CCD ID: MN) (formula: Mn).

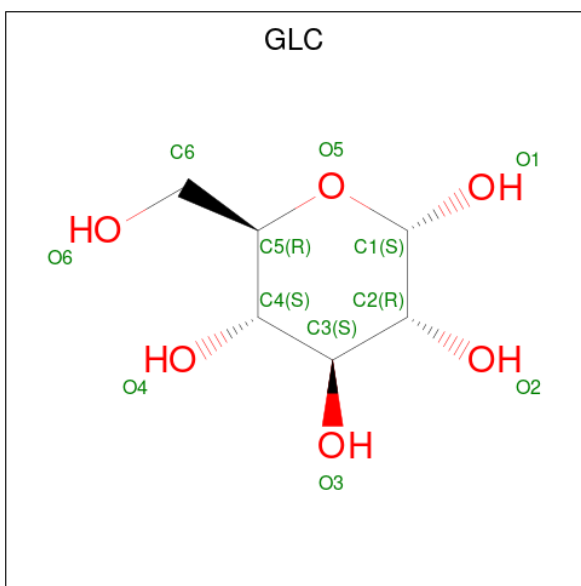
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Mn	0	0
			1	1		

- Molecule 3 is URIDINE-5'-DIPHOSPHATE (CCD ID: UDP) (formula: C₉H₁₄N₂O₁₂P₂).



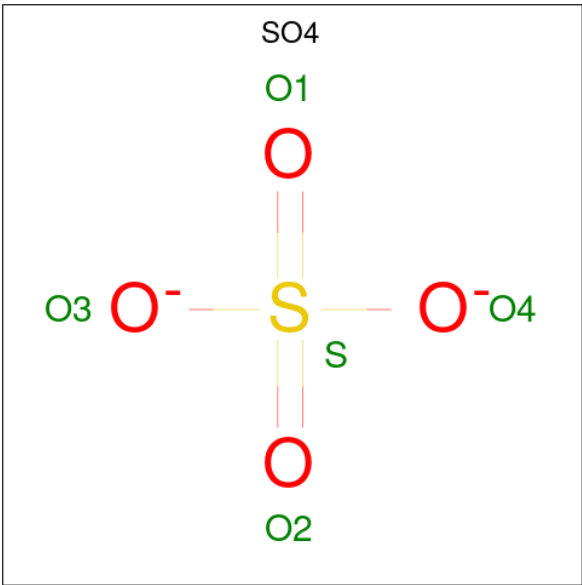
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			25	9	2	12	2		

- Molecule 4 is alpha-D-glucopyranose (CCD ID: GLC) (formula: $\text{C}_6\text{H}_{12}\text{O}_6$).



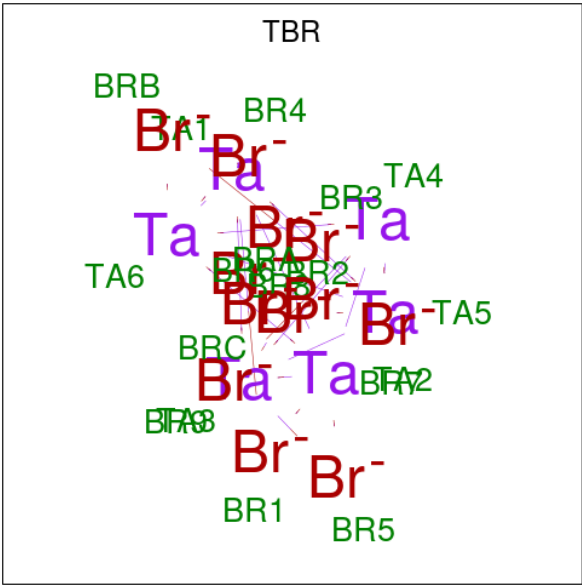
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			12	6	6		

- Molecule 5 is SULFATE ION (CCD ID: SO4) (formula: O_4S).



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

- Molecule 6 is HEXATANTALUM DODECABROMIDE (CCD ID: TBR) (formula: Br₁₂Ta₆).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	Br	Ta	0	0
			18	12	6		
6	A	1	Total	Br	Ta	0	0
			18	12	6		
6	A	1	Total	Br	Ta	0	0
			18	12	6		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	Br	Ta	0	0
			18	12	6		

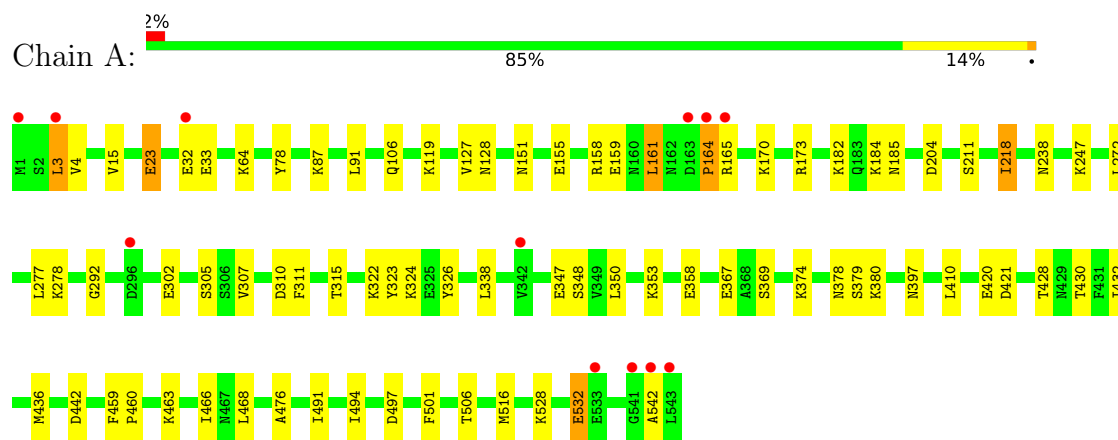
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	354	Total	O	0	0
			354	354		

3 Residue-property plots

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: TOXIN B



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	63.01Å 95.80Å 112.92Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.20 30.00 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.8 (30.00-2.20) 99.8 (30.00-2.20)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.62 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.197 , 0.244 0.199 , 0.247	Depositor DCC
R_{free} test set	1117 reflections (3.16%)	wwPDB-VP
Wilson B-factor (Å ²)	34.8	Xtriage
Anisotropy	0.190	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 41.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4902	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.41% 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: SO4, UDP, MN, TBR, GLC

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	0.69	0/4513	0.91	3/6089 (0.0%)

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	1

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	542	ALA	N-CA-C	-8.11	104.26	114.56
1	A	420	GLU	N-CA-C	-6.37	100.47	109.96
1	A	430	THR	N-CA-C	-5.25	105.64	111.36

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	164	PRO	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	4433	0	4320	49	0
2	A	1	0	0	0	0
3	A	25	0	11	0	0
4	A	12	0	12	0	0
5	A	5	0	0	0	0
6	A	72	0	0	16	0
7	A	354	0	0	3	0
All	All	4902	0	4343	49	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 6.

All (49) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:380:LYS:HG2	6:A:1549:TBR:BR3	2.25	0.91
1:A:516:MET:SD	6:A:1549:TBR:BRC	2.92	0.83
1:A:380:LYS:CG	6:A:1549:TBR:BR3	2.82	0.82
1:A:32:GLU:HG2	6:A:1548:TBR:BR7	2.37	0.79
1:A:378:ASN:ND2	6:A:1549:TBR:BRB	2.74	0.76
1:A:3:LEU:HD12	1:A:4:VAL:H	1.54	0.71
1:A:516:MET:HE1	6:A:1549:TBR:BRC	2.46	0.70
1:A:3:LEU:HG	1:A:78:TYR:HB3	1.74	0.69
1:A:170:LYS:HG3	1:A:173:ARG:HH21	1.57	0.69
1:A:323:TYR:CZ	1:A:347:GLU:HB3	2.30	0.66
1:A:428:THR:HG21	6:A:1551:TBR:BR3	2.52	0.65
1:A:32:GLU:CG	6:A:1548:TBR:BR7	3.01	0.64
1:A:204:ASP:OD1	7:A:2139:HOH:O	2.15	0.64
1:A:322:LYS:HE2	7:A:2202:HOH:O	1.99	0.63
1:A:528:LYS:O	1:A:532:GLU:HG2	1.99	0.62
1:A:151:ASN:O	1:A:155:GLU:HG3	1.99	0.62
1:A:311:PHE:O	1:A:315:THR:HG23	2.05	0.57
1:A:516:MET:CE	6:A:1549:TBR:BRC	3.09	0.55
1:A:494:ILE:HD11	6:A:1550:TBR:BR2	2.62	0.55
1:A:170:LYS:HG3	1:A:173:ARG:NH2	2.21	0.54
1:A:497:ASP:OD1	6:A:1550:TBR:BR2	2.82	0.53
1:A:459:PHE:CD2	1:A:460:PRO:HD2	2.45	0.52
1:A:379:SER:OG	6:A:1549:TBR:BR3	2.83	0.51
1:A:211:SER:HB2	1:A:218:ILE:HG22	1.92	0.51
1:A:410:LEU:HD22	1:A:468:LEU:HD21	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:158:ARG:O	1:A:161:LEU:HB2	2.12	0.49
1:A:23:GLU:H	1:A:23:GLU:CD	2.20	0.49
1:A:170:LYS:CG	1:A:173:ARG:HH21	2.25	0.48
1:A:353:LYS:HD2	1:A:358:GLU:HB3	1.94	0.48
1:A:33:GLU:HG3	6:A:1548:TBR:BRA	2.69	0.47
1:A:459:PHE:CG	1:A:460:PRO:HD2	2.50	0.47
1:A:278:LYS:HE3	1:A:397:ASN:OD1	2.16	0.46
1:A:374:LYS:HA	1:A:501:PHE:O	2.16	0.45
1:A:128:ASN:ND2	1:A:238:ASN:HD22	2.14	0.45
1:A:476:ALA:HB2	1:A:491:ILE:HD11	1.99	0.44
1:A:292:GLY:HA3	1:A:506:THR:O	2.19	0.43
1:A:3:LEU:HD13	1:A:3:LEU:HA	1.74	0.43
1:A:380:LYS:HG3	6:A:1549:TBR:BR3	2.69	0.43
1:A:463:LYS:HB3	1:A:466:ILE:HD12	2.01	0.42
1:A:310:ASP:OD2	6:A:1549:TBR:BR2	2.93	0.42
1:A:323:TYR:CE1	1:A:347:GLU:HB3	2.55	0.41
1:A:184:LYS:HD2	1:A:421:ASP:OD1	2.20	0.41
1:A:3:LEU:HD22	7:A:2055:HOH:O	2.20	0.41
1:A:324:LYS:HB3	1:A:326:TYR:CE1	2.56	0.41
1:A:87:LYS:O	1:A:369:SER:HB2	2.21	0.41
1:A:119:LYS:HG2	1:A:127:VAL:HG21	2.02	0.40
1:A:182:LYS:O	1:A:185:ASN:HB3	2.20	0.40
1:A:432:ILE:HD12	6:A:1551:TBR:BRB	2.76	0.40
1:A:436:MET:HE2	1:A:436:MET:HB3	1.96	0.40

There are no symmetry-related clashes.

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	541/543 (100%)	526 (97%)	14 (3%)	1 (0%)	43 51

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	164	PRO

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	495/495 (100%)	473 (96%)	22 (4%)	25	34

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	LEU
1	A	15	VAL
1	A	23	GLU
1	A	64	LYS
1	A	91	LEU
1	A	106	GLN
1	A	159	GLU
1	A	161	LEU
1	A	165	ARG
1	A	218	ILE
1	A	247	LYS
1	A	272	LEU
1	A	277	LEU
1	A	302	GLU
1	A	305	SER
1	A	307	VAL
1	A	338	LEU
1	A	348	SER
1	A	350	LEU
1	A	367	GLU
1	A	442	ASP
1	A	532	GLU

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	20	GLN
1	A	106	GLN
1	A	113	ASN
1	A	128	ASN
1	A	412	ASN
1	A	415	ASN
1	A	422	ASN
1	A	507	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 8 ligands modelled in this entry, 1 is monoatomic - leaving 7 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	TBR	A	1548	-	0,36,36	-	-	-		
4	GLC	A	1546	-	12,12,12	0.62	0	17,17,17	1.37	3 (17%)
3	UDP	A	1545	2	24,26,26	1.63	3 (12%)	37,40,40	1.82	8 (21%)
6	TBR	A	1550	-	0,36,36	-	-	-		
5	SO4	A	1547	-	4,4,4	0.20	0	6,6,6	0.24	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	TBR	A	1549	-	0,36,36	-	-	-		
6	TBR	A	1551	-	0,36,36	-	-	-		

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	GLC	A	1546	-	-	0/2/22/22	0/1/1/1
3	UDP	A	1545	2	-	0/16/32/32	0/2/2/2

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1545	UDP	PB-O1B	5.15	1.67	1.50
3	A	1545	UDP	O4'-C1'	2.56	1.48	1.42
3	A	1545	UDP	C2-N3	2.56	1.42	1.38

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1545	UDP	C4-N3-C2	-5.46	119.38	126.58
3	A	1545	UDP	C5-C4-N3	3.93	120.73	114.84
3	A	1545	UDP	N3-C2-N1	3.56	119.62	114.89
3	A	1545	UDP	O3B-PB-O3A	3.49	116.33	104.64
3	A	1545	UDP	PA-O3A-PB	-3.37	121.28	132.83
3	A	1545	UDP	O4-C4-C5	-3.29	119.38	125.16
3	A	1545	UDP	O2-C2-N1	-3.12	118.63	122.79
4	A	1546	GLC	O5-C1-C2	2.64	114.99	110.28
4	A	1546	GLC	O2-C2-C3	-2.37	104.86	110.35
3	A	1545	UDP	O4'-C4'-C3'	-2.07	101.01	105.11
4	A	1546	GLC	C1-C2-C3	2.05	114.56	110.31

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 16 short contacts:

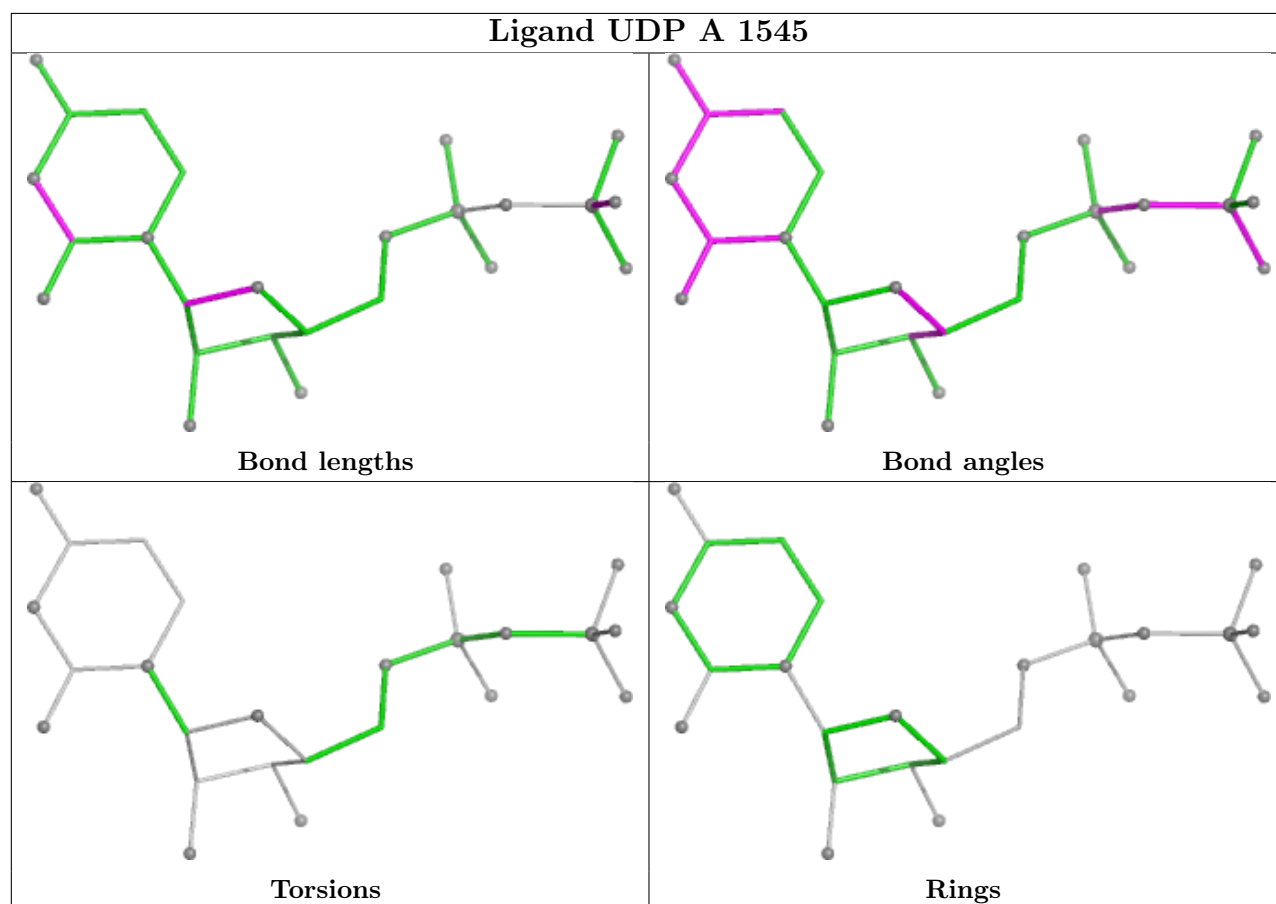
Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	1548	TBR	3	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	1550	TBR	2	0
6	A	1549	TBR	9	0
6	A	1551	TBR	2	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 [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	543/543 (100%)	0.31	12 (2%) 62 59	28, 40, 51, 64	0

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	163	ASP	5.7
1	A	164	PRO	5.5
1	A	543	LEU	3.7
1	A	542	ALA	3.5
1	A	533	GLU	3.2
1	A	541	GLY	3.2
1	A	32	GLU	2.9
1	A	296	ASP	2.7
1	A	3	LEU	2.7
1	A	342	VAL	2.3
1	A	165	ARG	2.3
1	A	1	MET	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 oligosaccharides in this entry.

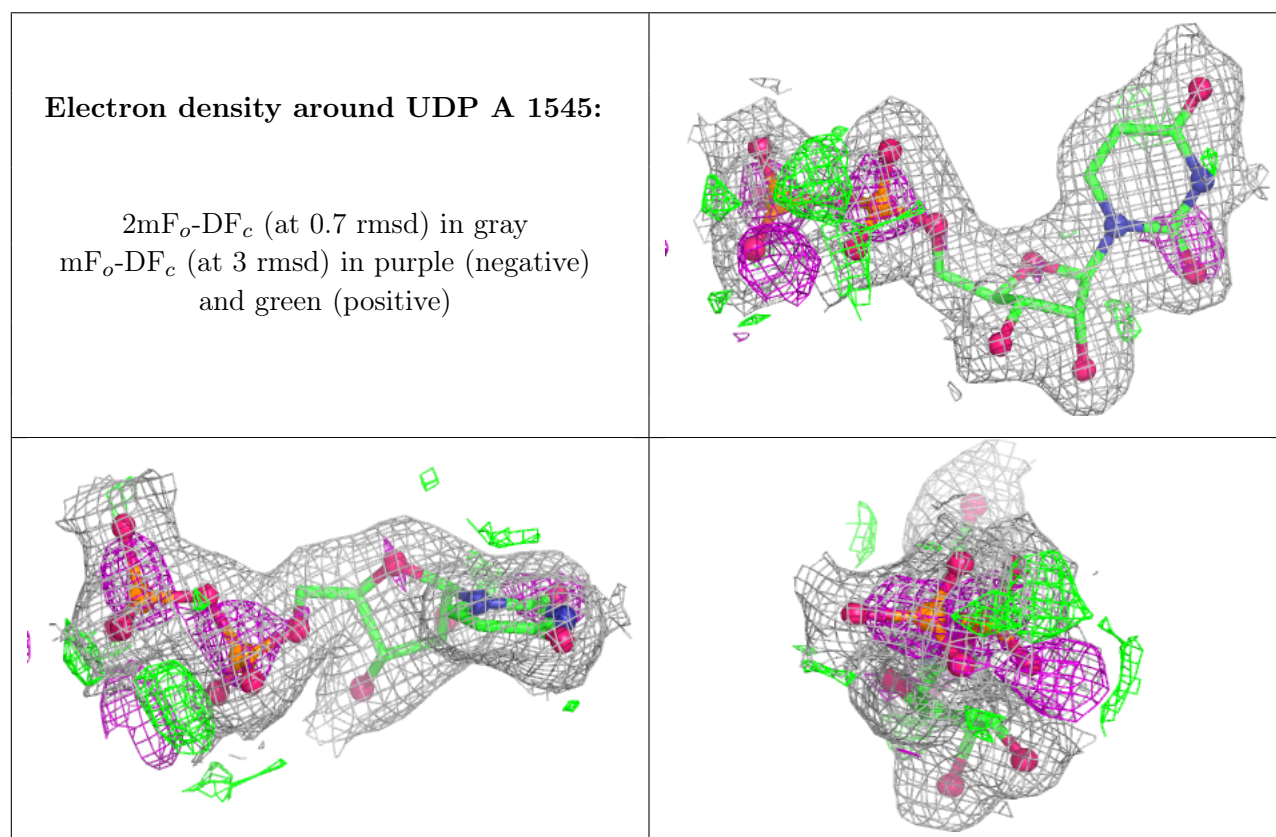
6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum,

median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
6	TBR	A	1550	18/18	0.77	0.22	55,56,56,58	18
6	TBR	A	1549	18/18	0.79	0.21	54,55,56,57	18
6	TBR	A	1551	18/18	0.83	0.19	51,52,52,53	18
6	TBR	A	1548	18/18	0.92	0.12	44,47,49,50	18
4	GLC	A	1546	12/12	0.96	0.09	23,25,28,30	0
5	SO4	A	1547	5/5	0.96	0.14	49,49,50,52	0
3	UDP	A	1545	25/25	0.96	0.08	22,26,29,29	0
2	MN	A	1544	1/1	0.99	0.08	25,25,25,25	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.



6.5 Other polymers ⓘ

There are no such residues in this entry.