



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 08:44 PM GMT

PDB ID : 1LTE
Title : STRUCTURE OF A LEGUME LECTIN WITH AN ORDERED N-LINKED CARBOHYDRATE IN COMPLEX WITH LACTOSE
Authors : Shaanan, B.; Lis, H.; Sharon, N.
Deposited on : 1991-06-25
Resolution : 2.00 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
<http://wwpdb.org/validation/2016/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 : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

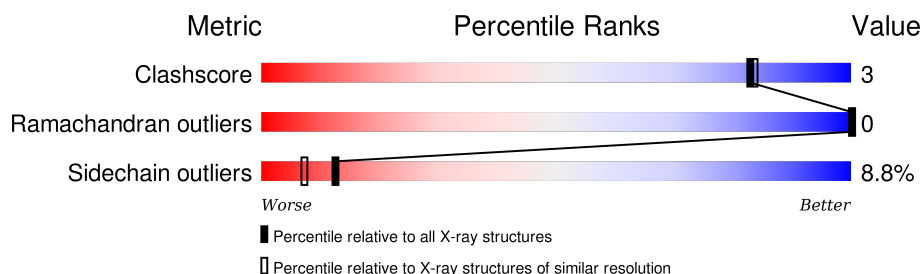
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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))
Clashscore	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (2.00-2.00)

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. 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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	239	 79% 17% . .

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 2062 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 CORAL TREE LECTIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	239	Total	C	N	O	S	0	0	0
			1857	1188	302	364	3			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	24	SER	ALA	CONFLICT	UNP P16404
A	114	GLN	ASN	CONFLICT	UNP P16404
A	178	LEU	ILE	CONFLICT	UNP P16404

- Molecule 2 is a polymer of unknown type called SUGAR (6-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	7	Total	C	N	O	0	0
			80	45	2	33		

- Molecule 3 is a polymer of unknown type called SUGAR (2-MER).

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	2	Total	C	O	0	0
			23	12	11		

- Molecule 4 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

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

- Molecule 5 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total 1	Ca 1	0	0

- Molecule 6 is water.

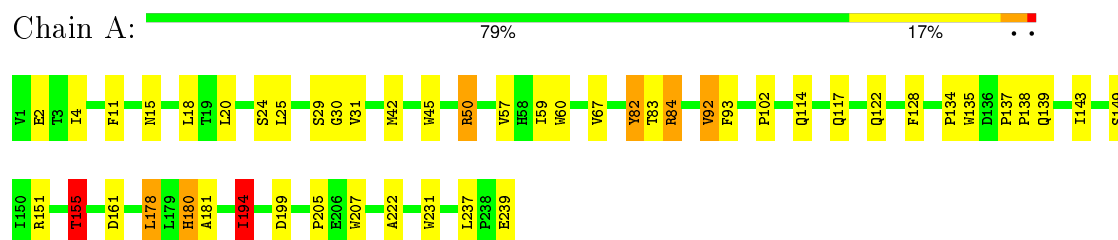
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	100	Total 100	O 100	0	0

3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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 ($\text{RSRZ} > 2$). 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.

Note EDS was not executed.

• Molecule 1: CORAL TREE LECTIN



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	84.40 Å 73.05 Å 71.40 Å 90.00° 113.42° 90.00°	Depositor
Resolution (Å)	(Not available) – 2.00	Depositor
% Data completeness (in resolution range)	(Not available) ((Not available)-2.00)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR	Depositor
R, R_{free}	0.190 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	2062	wwPDB-VP
Average B, all atoms (Å ²)	22.0	wwPDB-VP

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: BGC, NAG, CA, MN, GAL, BMA, MAN, XYL, FUC

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	0.77	0/1908	1.55	33/2606 (1.3%)

There are no bond length outliers.

All (33) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	50	ARG	NE-CZ-NH2	-15.25	112.67	120.30
1	A	50	ARG	NE-CZ-NH1	10.47	125.54	120.30
1	A	92	VAL	CB-CA-C	-8.81	94.66	111.40
1	A	199	ASP	CB-CG-OD1	8.70	126.13	118.30
1	A	60	TRP	CD1-CG-CD2	8.23	112.89	106.30
1	A	45	TRP	CD1-CG-CD2	7.71	112.47	106.30
1	A	135	TRP	CD1-CG-CD2	7.70	112.46	106.30
1	A	50	ARG	CB-CG-CD	-7.57	91.93	111.60
1	A	180	HIS	CA-CB-CG	7.51	126.36	113.60
1	A	60	TRP	CE2-CD2-CG	-7.30	101.46	107.30
1	A	207	TRP	CD1-CG-CD2	7.22	112.08	106.30
1	A	231	TRP	CD1-CG-CD2	6.97	111.88	106.30
1	A	45	TRP	CE2-CD2-CG	-6.95	101.74	107.30
1	A	207	TRP	CE2-CD2-CG	-6.77	101.88	107.30
1	A	194	ILE	CB-CA-C	-6.76	98.07	111.60
1	A	135	TRP	CE2-CD2-CG	-6.61	102.01	107.30
1	A	231	TRP	CE2-CD2-CG	-6.47	102.12	107.30
1	A	151	ARG	N-CA-CB	-6.13	99.56	110.60
1	A	2	GLU	CA-CB-CG	6.11	126.84	113.40
1	A	161	ASP	N-CA-C	-6.05	94.66	111.00
1	A	82	TYR	CB-CG-CD1	-5.95	117.43	121.00
1	A	25	LEU	CA-CB-CG	5.60	128.18	115.30
1	A	93	PHE	N-CA-C	-5.53	96.06	111.00
1	A	135	TRP	CG-CD1-NE1	-5.47	104.62	110.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	84	ARG	CG-CD-NE	-5.34	100.58	111.80
1	A	222	ALA	N-CA-C	-5.32	96.65	111.00
1	A	178	LEU	N-CA-CB	-5.23	99.94	110.40
1	A	45	TRP	CG-CD1-NE1	-5.19	104.91	110.10
1	A	42	MET	CA-CB-CG	-5.18	104.50	113.30
1	A	128	PHE	N-CA-C	-5.17	97.05	111.00
1	A	60	TRP	CG-CD1-NE1	-5.16	104.94	110.10
1	A	155	THR	N-CA-CB	-5.05	100.70	110.30
1	A	60	TRP	CA-CB-CG	5.02	123.23	113.70

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

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	1857	0	1789	12	0
2	A	80	0	68	0	0
3	A	23	0	21	0	0
4	A	1	0	0	0	0
5	A	1	0	0	0	0
6	A	100	0	0	1	0
All	All	2062	0	1878	12	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 3.

All (12) 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:134:PRO:HD3	6:A:814:HOH:O	1.96	0.64
1:A:137:PRO:HG3	1:A:155:THR:HG21	1.90	0.54
1:A:50:ARG:HD2	1:A:102:PRO:HG3	1.92	0.52
1:A:4:ILE:HD13	1:A:57:VAL:HG12	1.90	0.52
1:A:122:GLN:HB3	1:A:205:PRO:HD3	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:117:GLN:HG3	1:A:149:SER:HB2	1.94	0.50
1:A:181:ALA:HB3	1:A:194:ILE:CD1	2.42	0.50
1:A:138:PRO:HD2	1:A:139:GLN:OE1	2.13	0.48
1:A:181:ALA:HB3	1:A:194:ILE:HD13	1.95	0.47
1:A:11:PHE:O	1:A:30:GLY:HA2	2.15	0.46
1:A:59:ILE:HA	1:A:59:ILE:HD12	1.92	0.42
1:A:143:ILE:HG21	1:A:194:ILE:HD11	2.04	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	237/239 (99%)	232 (98%)	5 (2%)	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	205/205 (100%)	187 (91%)	18 (9%)	12	7

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	15	ASN
1	A	18	LEU
1	A	20	LEU
1	A	24	SER
1	A	29	SER
1	A	31	VAL
1	A	67	VAL
1	A	82	TYR
1	A	83	THR
1	A	84	ARG
1	A	92	VAL
1	A	114	GLN
1	A	155	THR
1	A	178	LEU
1	A	180	HIS
1	A	194	ILE
1	A	237	LEU
1	A	239	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 ⓘ

9 carbohydrates 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	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	NAG	A	301	1,2	14,14,15	0.67	0	15,19,21	1.63	3 (20%)
2	FUC	A	302	2	10,10,11	0.55	0	14,14,16	1.10	1 (7%)
2	NAG	A	303	2	14,14,15	0.48	0	15,19,21	1.09	2 (13%)
2	BMA	A	304	2	11,11,12	0.53	0	14,15,17	0.94	0
2	XYL	A	305	2	7,8,9	0.58	0	7,10,11	1.16	1 (14%)
2	MAN	A	306	2	11,11,12	0.78	0	14,15,17	0.93	0
2	MAN	A	307	2	11,11,12	0.58	0	14,15,17	1.08	1 (7%)
3	BGC	A	401	3	12,12,12	0.83	0	17,17,17	1.25	1 (5%)
3	GAL	A	402	3	11,11,12	0.67	0	14,15,17	1.31	3 (21%)

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	NAG	A	301	1,2	-	0/6/23/26	0/1/1/1
2	FUC	A	302	2	-	0/0/17/20	0/1/1/1
2	NAG	A	303	2	-	0/6/23/26	0/1/1/1
2	BMA	A	304	2	-	0/2/19/22	0/1/1/1
2	XYL	A	305	2	-	0/10/10/12	0/0/0/0
2	MAN	A	306	2	-	0/2/19/22	0/1/1/1
2	MAN	A	307	2	-	0/2/19/22	0/1/1/1
3	BGC	A	401	3	-	0/2/22/22	0/1/1/1
3	GAL	A	402	3	-	0/2/19/22	0/1/1/1

There are no bond length outliers.

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	305	XYL	C4-C3-C2	-2.40	108.56	112.47
2	A	301	NAG	O3-C3-C2	-2.24	104.67	109.11
3	A	402	GAL	C2-C3-C4	-2.24	107.23	111.04
2	A	303	NAG	C3-C2-N2	-2.18	105.35	110.56
2	A	303	NAG	C3-C4-C5	2.23	114.09	110.20
2	A	307	MAN	C1-O5-C5	2.25	115.10	112.25
3	A	402	GAL	O2-C2-C1	2.33	113.89	109.21
2	A	301	NAG	C8-C7-N2	2.34	120.59	116.11
2	A	302	FUC	C1-O5-C5	2.44	116.15	112.38
3	A	401	BGC	C1-O5-C5	2.85	118.74	113.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	402	GAL	C1-O5-C5	2.98	116.04	112.25
2	A	301	NAG	C1-O5-C5	4.50	117.96	112.25

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.6 Ligand geometry [i](#)

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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 ⓘ

6.1 Protein, DNA and RNA chains ⓘ

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates ⓘ

EDS was not executed - this section will therefore be empty.

6.4 Ligands ⓘ

EDS was not executed - this section will therefore be empty.

6.5 Other polymers ⓘ

EDS was not executed - this section will therefore be empty.