WORLDWIDE PDB PROTEIN DATA BANK

wwPDB X-ray Structure Validation Summary Report (i)

Jan 31, 2016 – 08:39 PM GMT

PDB ID	:	1LBI
Title	:	LAC REPRESSOR
Authors	:	Lewis, M.; Chang, G.; Horton, N.C.; Kercher, M.A.; Pace, H.C.; Lu, P.
Deposited on	:	1996-02-17
Resolution	:	2.70 Å(reported)

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

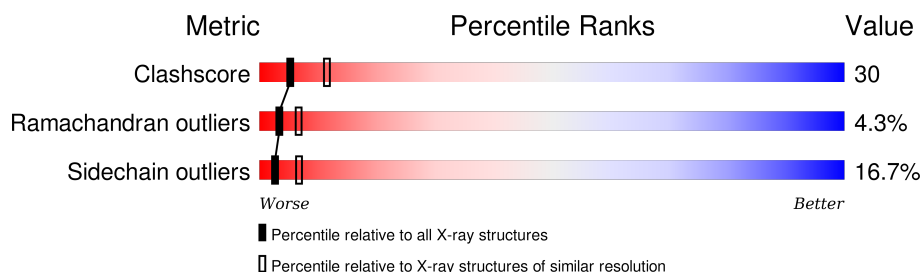|		
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

1 Overall quality at a glance (i)

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

The reported resolution of this entry is 2.70 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	Percentile Ranks	Value
Clashscore		30
Ramachandran outliers		4.3%
Sidechain outliers		16.7%

Worse   *Better*

■ Percentile relative to all X-ray structures

▯ Percentile relative to X-ray structures of similar resolution

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	102246	2422 (2.70-2.70)
Ramachandran outliers	100387	2382 (2.70-2.70)
Sidechain outliers	100360	2382 (2.70-2.70)

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 <=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	360	32% 41% 9% 18%
1	B	360	37% 35% 10% • 18%
1	C	360	37% 36% 9% • 18%
1	D	360	34% 39% 8% • 18%

2 Entry composition (i)

There is only 1 type of molecule in this entry. The entry contains 8872 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 LAC REPRESSOR.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	296	Total 2218	C 1383	N 396	O 428	S 11	0	0	0
1	B	296	Total 2218	C 1383	N 396	O 428	S 11	0	0	0
1	C	296	Total 2218	C 1383	N 396	O 428	S 11	0	0	0
1	D	296	Total 2218	C 1383	N 396	O 428	S 11	0	0	0

There are 8 discrepancies between the modelled and reference sequences:

| Chain	Residue	Modelled	Actual	Comment	Reference
A	109	THR	ALA	CONFLICT	UNP P03023
A	286	LEU	SER	CONFLICT	UNP P03023
B	109	THR	ALA	CONFLICT	UNP P03023
B	286	LEU	SER	CONFLICT	UNP P03023
C	109	THR	ALA	CONFLICT	UNP P03023
C	286	LEU	SER	CONFLICT	UNP P03023
D	109	THR	ALA	CONFLICT	UNP P03023
D	286	LEU	SER	CONFLICT	UNP P03023

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 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 (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.

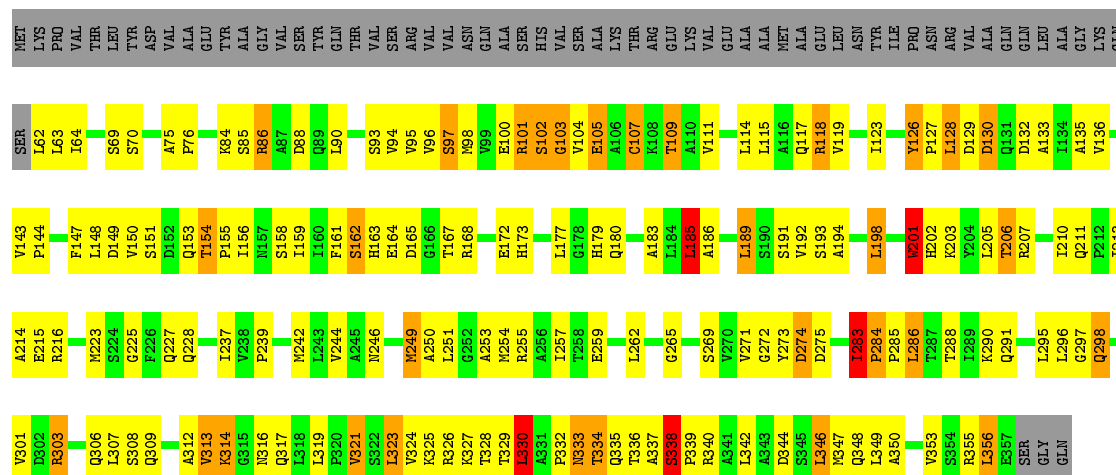Note EDS was not executed.

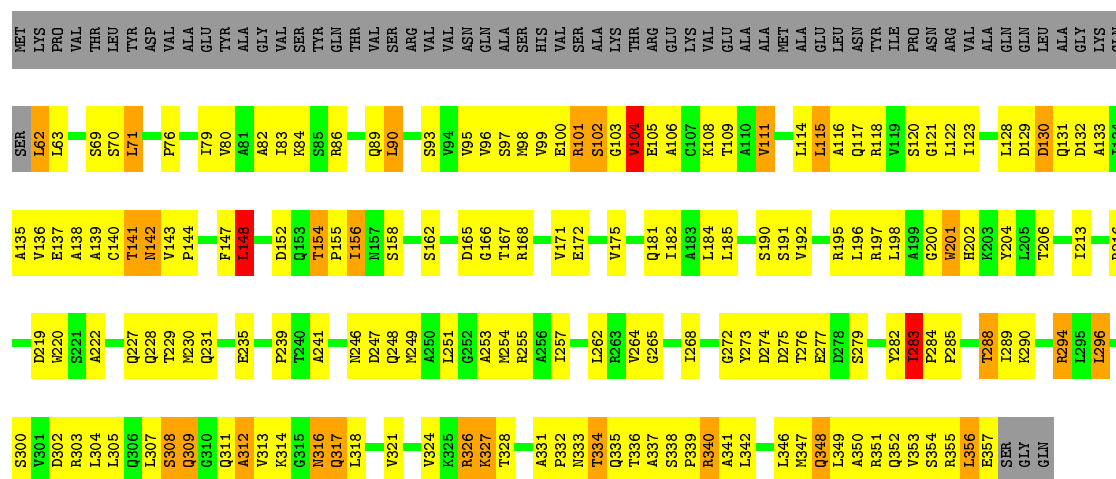- Molecule 1: LAC REPRESSOR
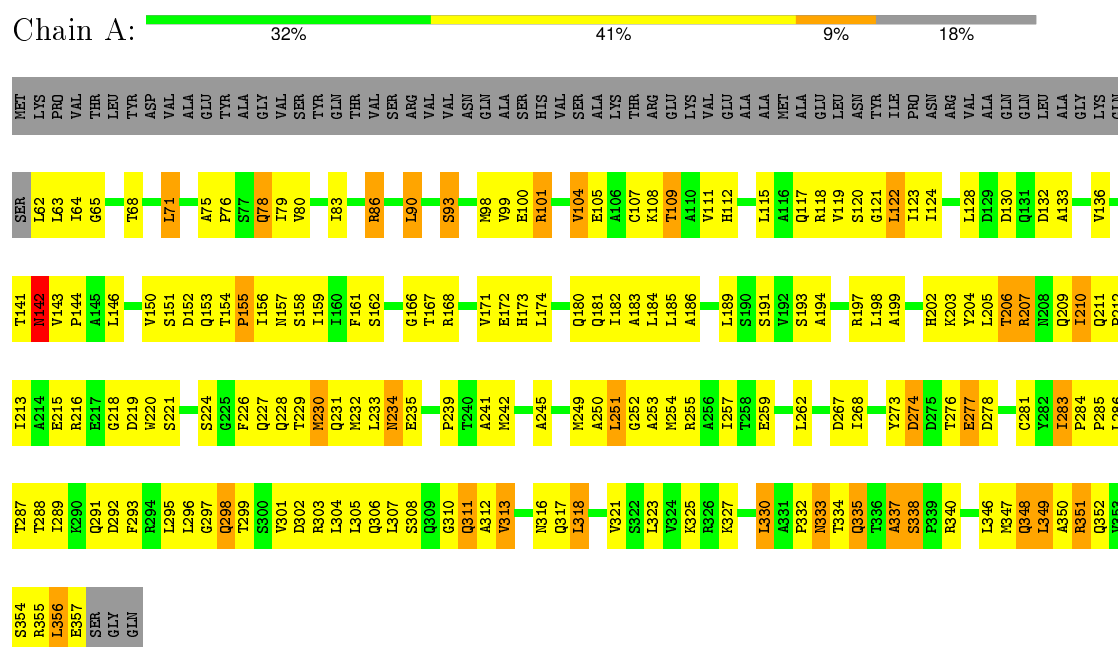
- Molecule 1: LAC REPRESSOR
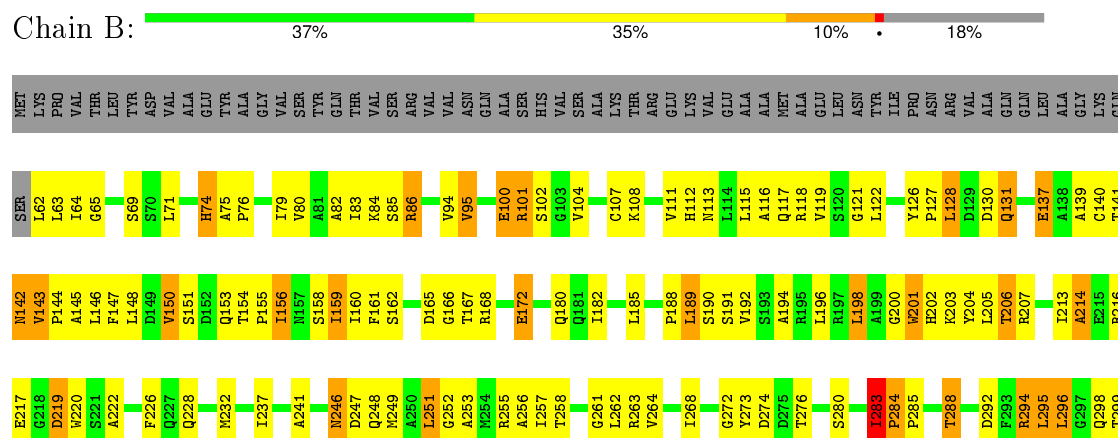
• Molecule 1: LAC REPRESSOR

Chain C:

| 37% | 36% | 9% | • | 18% |

• Molecule 1: LAC REPRESSOR

Chain D:

| 34% | 39% | 8% | • | 18% |

4 Data and refinement statistics (i)

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	160.30Å 73.30Å 147.80Å 90.00° 120.30° 90.00°	Depositor
Resolution (Å)	15.00 – 2.70	Depositor
% Data completeness (in resolution range)	(Not available) (15.00-2.70)	Depositor
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.250 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	8872	wwPDB-VP
Average B, all atoms (Å^2)	25.0	wwPDB-VP

5 Model quality (i)

5.1 Standard geometry (i)

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	0.78	0/2247	1.04	4/3055 (0.1%)
1	B	0.82	0/2247	1.11	10/3055 (0.3%)
1	C	0.84	1/2247 (0.0%)	1.04	9/3055 (0.3%)
1	D	0.85	0/2247	0.99	3/3055 (0.1%)
All	All	0.82	1/8988 (0.0%)	1.05	26/12220 (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	C	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	201	TRP	CB-CG	6.07	1.61	1.50

The worst 5 of 26 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	283	ILE	C-N-CD	-19.64	77.40	120.60
1	B	283	ILE	C-N-CA	10.29	165.20	122.00
1	C	330	LEU	CA-CB-CG	7.59	132.76	115.30
1	C	283	ILE	C-N-CD	7.05	143.21	128.40
1	B	336	THR	N-CA-C	-6.62	93.13	111.00

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	126	TYR	Sidechain

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	2218	0	2269	158	0
1	B	2218	0	2269	140	0
1	C	2218	0	2269	148	0
1	D	2218	0	2269	145	0
All	All	8872	0	9076	544	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 30.

The worst 5 of 544 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:144:PRO:HG2	1:A:308:SER:HA	1.24	1.13
1:A:115:LEU:HD11	1:A:143:VAL:HG11	1.18	1.10
1:C:283:ILE:HG22	1:C:284:PRO:HD3	1.43	1.01
1:B:189:LEU:HD13	1:B:217:GLU:HG3	1.40	0.99
1:A:144:PRO:HB3	1:A:307:LEU:HD22	1.44	0.98

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	294/360 (82%)	244 (83%)	32 (11%)	18 (6%)	2	3
1	B	294/360 (82%)	247 (84%)	31 (10%)	16 (5%)	2	4
1	C	294/360 (82%)	260 (88%)	26 (9%)	8 (3%)	6	16
1	D	294/360 (82%)	247 (84%)	38 (13%)	9 (3%)	5	12
All	All	1176/1440 (82%)	998 (85%)	127 (11%)	51 (4%)	3	7

5 of 51 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	141	THR
1	A	142	ASN
1	A	162	SER
1	A	274	ASP
1	A	277	GLU

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	243/295 (82%)	208 (86%)	35 (14%)	4	10
1	B	243/295 (82%)	203 (84%)	40 (16%)	3	7
1	C	243/295 (82%)	201 (83%)	42 (17%)	2	6
1	D	243/295 (82%)	198 (82%)	45 (18%)	2	5
All	All	972/1180 (82%)	810 (83%)	162 (17%)	3	7

5 of 162 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	351	ARG
1	C	154	THR
1	D	313	VAL
1	C	69	SER
1	C	107	CYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 32 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	228	GLN
1	C	181	GLN
1	D	309	GLN
1	C	163	HIS
1	C	202	HIS

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 carbohydrates in this entry.

5.6 Ligand geometry (i)

There are no ligands in this entry.

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)

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

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

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

6.3 Carbohydrates (i)

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

6.4 Ligands (i)

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

6.5 Other polymers (i)

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