



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 10:50 PM GMT

PDB ID : 1VFD
Title : HUMAN LACTOFERRIN, N-TERMINAL LOBE MUTANT WITH ARG 121
REPLACED BY GLU (R121E)
Authors : Faber, H.R.; Day, C.L.; Baker, E.N.
Deposited on : 1996-10-01
Resolution : 2.50 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) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

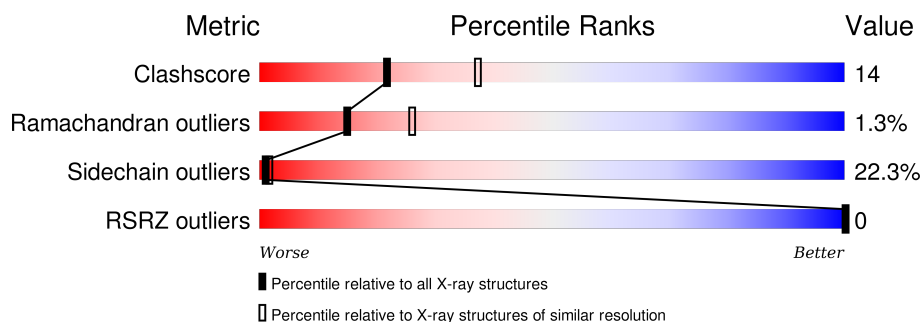
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	4242 (2.50-2.50)
Ramachandran outliers	100387	4156 (2.50-2.50)
Sidechain outliers	100360	4158 (2.50-2.50)
RSRZ outliers	91569	3562 (2.50-2.50)

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.

Mol	Chain	Length	Quality of chain
1	A	330	

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 2589 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 LACTOFERRIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	321	2500	1581	444	462	13	0	0	0

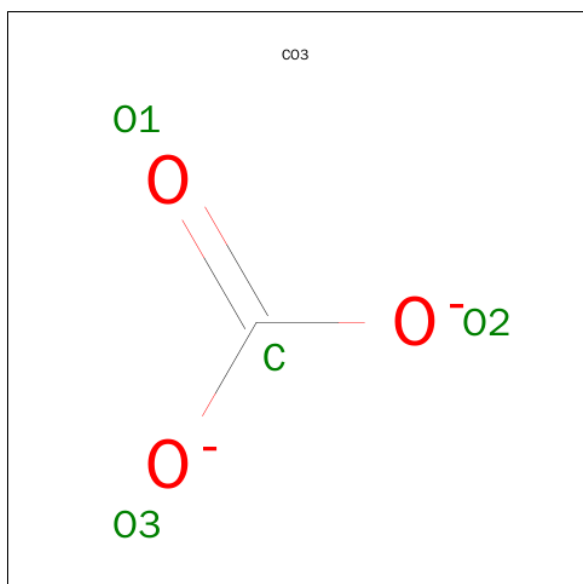
There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	28	ARG	LYS	VARIANT	UNP P02788
A	121	GLU	ARG	ENGINEERED	UNP P02788

- Molecule 2 is FE (III) ION (three-letter code: FE) (formula: Fe).

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

- Molecule 3 is CARBONATE ION (three-letter code: CO3) (formula: CO₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			4	1	3		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	84	Total	O	0	0
			84	84		

i

- Molecule 1: LACTOFERRIN



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	98.10 Å 78.90 Å 113.60 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.50 32.40 – 2.50	Depositor EDS
% Data completeness (in resolution range)	85.4 (20.00-2.50) 83.8 (32.40-2.50)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.65 (at 2.51 Å)	Xtriage
Refinement program	TNT	Depositor
R, R_{free}	0.204 , 0.310 0.184 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	44.3	Xtriage
Anisotropy	0.097	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.20 , 77.8	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.28$	Xtriage
Outliers	0 of 13063 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	2589	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.63% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: CO3, FE

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	0.84	2/2564 (0.1%)	1.16	10/3471 (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	1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	314	ILE	C-N	-5.86	1.20	1.34
1	A	192	TYR	CD2-CE2	-5.56	1.31	1.39

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	312	PRO	O-C-N	7.85	135.27	122.70
1	A	313	ARG	NE-CZ-NH2	7.51	124.06	120.30
1	A	224	ARG	NE-CZ-NH1	-6.75	116.92	120.30
1	A	55	ASP	N-CA-C	6.63	128.89	111.00
1	A	109	LEU	N-CA-C	6.56	128.72	111.00
1	A	314	ILE	O-C-N	-6.48	112.33	122.70
1	A	249	ARG	N-CA-C	-5.35	96.55	111.00
1	A	288	LEU	CB-CG-CD2	-5.18	102.19	111.00
1	A	301	LYS	CB-CA-C	-5.16	100.08	110.40
1	A	312	PRO	CA-C-N	-5.01	106.18	117.20

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	314	ILE	Mainchain

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	2500	0	2428	69	0
2	A	1	0	0	0	0
3	A	4	0	0	0	0
4	A	84	0	0	0	0
All	All	2589	0	2428	69	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 14.

All (69) 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:106:LEU:HD23	1:A:232:PRO:HA	1.70	0.72
1:A:61:GLY:HA2	1:A:64:ILE:HD12	1.74	0.69
1:A:168:ASN:HA	1:A:171:ARG:HB2	1.75	0.69
1:A:133:ARG:HA	1:A:136:LEU:HB2	1.75	0.68
1:A:251:PRO:HG3	1:A:320:LEU:HA	1.77	0.67
1:A:116:HIS:CD2	1:A:158:VAL:HG22	2.29	0.66
1:A:95:VAL:HB	1:A:229:LEU:HD22	1.78	0.64
1:A:136:LEU:HD12	1:A:145:ILE:HD11	1.79	0.63
1:A:26:MET:HE3	1:A:32:PRO:HD2	1.79	0.63
1:A:119:LEU:HG	1:A:120:ARG:HG3	1.80	0.63
1:A:104:PHE:O	1:A:236:ARG:NH1	2.33	0.62
1:A:122:THR:HG21	1:A:250:VAL:HG11	1.82	0.62
1:A:258:ARG:HB2	1:A:262:GLY:HA2	1.81	0.61
1:A:155:ALA:HB1	1:A:171:ARG:HB3	1.82	0.61
1:A:155:ALA:HA	1:A:168:ASN:HB2	1.83	0.60
1:A:49:ILE:HG12	1:A:55:ASP:HA	1.85	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:113:LYS:HB2	1:A:204:GLY:HA2	1.88	0.54
1:A:109:LEU:HD13	1:A:153:PHE:HE1	1.72	0.53
1:A:12:SER:HA	1:A:38:LYS:HE3	1.90	0.52
1:A:21:GLN:O	1:A:25:ASN:ND2	2.42	0.52
1:A:48:ALA:HA	1:A:51:GLU:HG2	1.92	0.52
1:A:187:GLU:HB3	1:A:190:PHE:HB2	1.92	0.52
1:A:157:CYS:HB2	1:A:172:LEU:HB2	1.91	0.51
1:A:18:LYS:HG2	1:A:299:LEU:HD12	1.92	0.51
1:A:75:ARG:NH1	1:A:264:GLU:OE2	2.44	0.51
1:A:51:GLU:HG3	1:A:53:ARG:HG3	1.94	0.50
1:A:7:GLN:HA	1:A:35:SER:HB3	1.94	0.50
1:A:59:LEU:HD12	1:A:256:VAL:HG11	1.94	0.50
1:A:49:ILE:HA	1:A:54:ALA:O	2.11	0.50
1:A:114:SER:HB2	1:A:208:PHE:HE2	1.77	0.49
1:A:136:LEU:O	1:A:137:ASN:HB3	2.13	0.49
1:A:231:CYS:SG	1:A:237:LYS:HD2	2.53	0.49
1:A:191:SER:OG	1:A:192:TYR:N	2.45	0.49
1:A:74:LEU:HB3	1:A:256:VAL:HG22	1.95	0.48
1:A:196:PHE:CZ	1:A:214:VAL:HG13	2.49	0.48
1:A:209:ILE:HD12	1:A:213:THR:HB	1.95	0.48
1:A:189:TYR:CD2	1:A:198:CYS:HB2	2.49	0.48
1:A:313:ARG:HA	1:A:313:ARG:HD2	1.75	0.47
1:A:97:VAL:HG21	1:A:209:ILE:HD11	1.96	0.47
1:A:118:GLY:HA2	1:A:159:PRO:HG2	1.95	0.46
1:A:291:SER:HB3	1:A:298:LEU:HD22	1.98	0.46
1:A:128:PRO:HG3	1:A:247:LEU:HD12	1.97	0.46
1:A:125:TRP:CZ3	1:A:158:VAL:HG11	2.51	0.46
1:A:25:ASN:HB3	1:A:278:PHE:CZ	2.52	0.45
1:A:224:ARG:HD3	1:A:224:ARG:HH11	1.61	0.45
1:A:189:TYR:HE1	1:A:197:LYS:HG2	1.82	0.45
1:A:192:TYR:CE1	1:A:210:ARG:HG2	2.51	0.44
1:A:81:VAL:HG23	1:A:307:PHE:HA	1.99	0.44
1:A:278:PHE:CD1	1:A:286:PHE:HB3	2.52	0.44
1:A:107:ASN:HD22	1:A:107:ASN:H	1.65	0.44
1:A:39:ARG:HG2	1:A:44:GLN:HB3	2.00	0.43
1:A:189:TYR:CE1	1:A:197:LYS:HG2	2.53	0.43
1:A:75:ARG:HD2	1:A:313:ARG:HD2	1.99	0.43
1:A:11:VAL:HB	1:A:15:GLU:OE1	2.18	0.43
1:A:26:MET:SD	1:A:278:PHE:HE2	2.42	0.43
1:A:105:GLN:HB2	1:A:107:ASN:ND2	2.33	0.43
1:A:209:ILE:HD12	1:A:213:THR:CB	2.48	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:118:GLY:HA3	1:A:121:GLU:HG3	2.00	0.42
1:A:32:PRO:HD3	1:A:273:GLN:CD	2.39	0.42
1:A:269:ASN:OD1	1:A:272:ARG:NH1	2.53	0.42
1:A:165:GLN:HB2	1:A:166:PHE:CE2	2.55	0.41
1:A:37:ILE:HD13	1:A:37:ILE:HA	1.93	0.41
1:A:113:LYS:O	1:A:205:ASP:N	2.45	0.41
1:A:209:ILE:HB	1:A:210:ARG:H	1.71	0.41
1:A:113:LYS:HB3	1:A:172:LEU:HD11	2.02	0.41
1:A:70:ALA:HA	1:A:71:PRO:HA	1.75	0.40
1:A:18:LYS:HG2	1:A:299:LEU:CD1	2.51	0.40
1:A:117:THR:HB	1:A:191:SER:HA	2.03	0.40
1:A:240:ASP:N	1:A:240:ASP:OD1	2.55	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	319/330 (97%)	293 (92%)	22 (7%)	4 (1%)	15	26

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	109	LEU
1	A	209	ILE
1	A	221	GLU
1	A	264	GLU

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	264/271 (97%)	205 (78%)	59 (22%)	1 2

All (59) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	SER
1	A	18	LYS
1	A	23	GLN
1	A	24	ARG
1	A	26	MET
1	A	29	VAL
1	A	39	ARG
1	A	49	ILE
1	A	52	ASN
1	A	75	ARG
1	A	85	GLU
1	A	86	ARG
1	A	87	GLN
1	A	89	ARG
1	A	97	VAL
1	A	99	LYS
1	A	105	GLN
1	A	107	ASN
1	A	109	LEU
1	A	113	LYS
1	A	136	LEU
1	A	143	GLU
1	A	146	GLU
1	A	154	SER
1	A	162	ASP
1	A	163	LYS
1	A	176	THR
1	A	178	GLU
1	A	187	GLU
1	A	191	SER

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Mol	Chain	Res	Type
1	A	197	LYS
1	A	198	CYS
1	A	210	ARG
1	A	211	GLU
1	A	214	VAL
1	A	218	LEU
1	A	219	SER
1	A	220	ASP
1	A	224	ARG
1	A	233	ASP
1	A	240	ASP
1	A	241	LYS
1	A	249	ARG
1	A	250	VAL
1	A	252	SER
1	A	261	ASN
1	A	263	LYS
1	A	265	ASP
1	A	276	GLU
1	A	280	LYS
1	A	291	SER
1	A	293	SER
1	A	296	LYS
1	A	298	LEU
1	A	301	LYS
1	A	303	SER
1	A	308	SER
1	A	314	ILE
1	A	322	SER

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

Mol	Chain	Res	Type
1	A	13	GLN
1	A	23	GLN
1	A	25	ASN
1	A	87	GLN
1	A	91	HIS
1	A	105	GLN
1	A	107	ASN
1	A	273	GLN

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 2 ligands modelled in this entry, 1 is monoatomic - leaving 1 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	$\# Z > 2$
3	CO3	A	401	2	0,3,3	0.00	-	0,3,3	0.00	-

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
3	CO3	A	401	2	-	0/0/0/0	0/0/0/0

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

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	321/330 (97%)	-0.73	0 100 100	13, 31, 66, 78	0

There are no RSRZ outliers to report.

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
3	CO3	A	401	4/4	0.96	0.20	1.04	22,26,31,33	0
2	FE	A	400	1/1	1.00	0.13	-2.12	23,23,23,23	0

6.5 Other polymers [i](#)

There are no such residues in this entry.