



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 04:01 PM UTC

PDB ID : 1VFD / pdb_00001vfd
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

<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	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

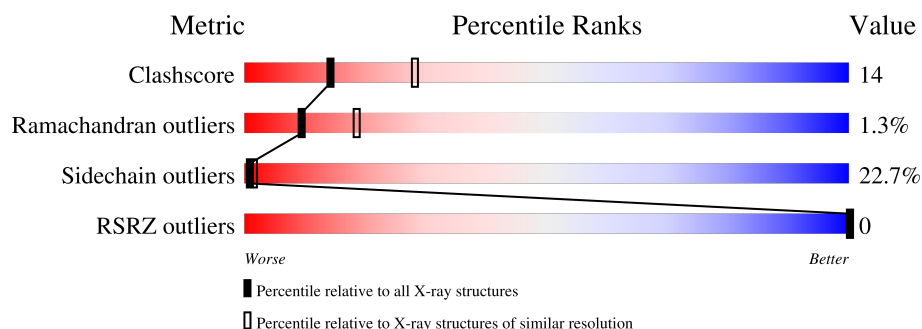
1 Overall quality at a glance

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	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (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 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	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

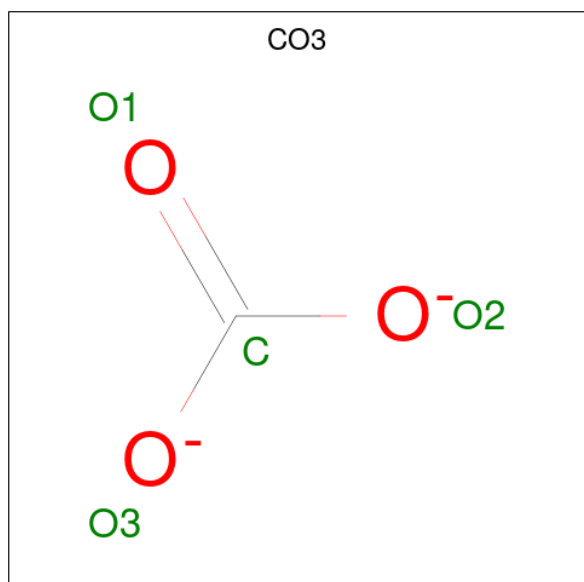
There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	28	ARG	LYS	variant	UNP P02788
A	121	GLU	ARG	engineered mutation	UNP P02788

- Molecule 2 is FE (III) ION (CCD ID: FE) (formula: Fe).

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

- Molecule 3 is CARBONATE ION (CCD ID: CO3) (formula: CO₃).



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

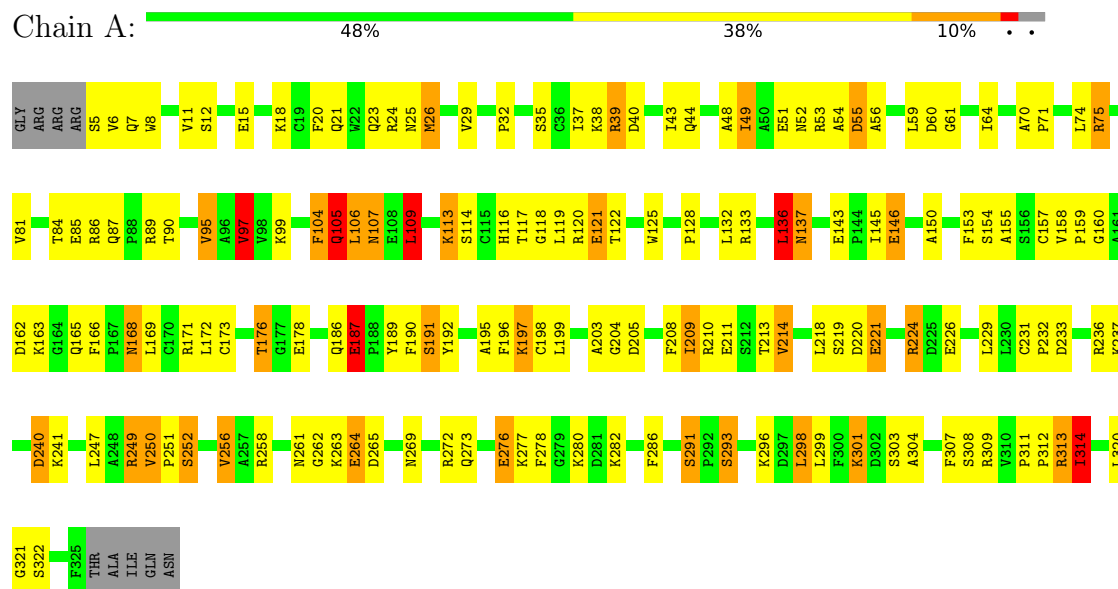
- Molecule 4 is water.

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

3 Residue-property plots [i](#)

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: LACTOFERRIN



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	98.10Å 78.90Å 113.60Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.50 20.00 – 2.50	Depositor EDS
% Data completeness (in resolution range)	85.4 (20.00-2.50) 83.6 (20.00-2.50)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.65 (at 2.51Å)	Xtriage
Refinement program	TNT	Depositor
R, R_{free}	0.204 , 0.310 0.186 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	44.3	Xtriage
Anisotropy	0.097	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.18 , 59.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	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, 0.333 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: FE, CO3

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	1.08	7/2564 (0.3%)	1.62	45/3471 (1.3%)

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 (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	314	ILE	C-N	-10.26	1.20	1.33
1	A	97	VAL	CA-CB	-7.17	1.45	1.54
1	A	195	ALA	CA-CB	-6.30	1.42	1.53
1	A	137	ASN	N-CA	6.05	1.53	1.46
1	A	95	VAL	CA-CB	-5.98	1.46	1.54
1	A	309	ARG	C-N	-5.54	1.27	1.33
1	A	95	VAL	CA-C	-5.19	1.47	1.53

All (45) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	55	ASP	N-CA-C	10.63	128.89	114.12
1	A	136	LEU	N-CA-C	9.80	123.19	111.82
1	A	312	PRO	O-C-N	9.75	135.27	122.89
1	A	143	GLU	N-CA-C	-9.11	98.63	109.93
1	A	311	PRO	O-C-N	8.49	125.22	121.31
1	A	109	LEU	N-CA-C	8.41	128.72	110.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	187	GLU	N-CA-C	-8.21	97.83	109.24
1	A	249	ARG	N-CA-C	-8.10	96.55	109.59
1	A	169	LEU	N-CA-C	-7.99	103.53	113.28
1	A	277	LYS	N-CA-C	7.74	121.21	111.69
1	A	314	ILE	O-C-N	-7.38	112.33	122.59
1	A	203	ALA	N-CA-C	7.16	118.88	111.14
1	A	293	SER	N-CA-C	7.05	120.53	109.96
1	A	321	GLY	N-CA-C	-6.89	104.58	112.29
1	A	276	GLU	N-CA-C	-6.60	103.57	111.69
1	A	309	ARG	N-CA-C	6.48	119.17	110.35
1	A	20	PHE	N-CA-C	-6.39	104.97	112.89
1	A	304	ALA	N-CA-C	-6.03	100.46	110.17
1	A	301	LYS	CB-CA-C	-5.99	100.08	109.84
1	A	40	ASP	N-CA-C	5.92	120.66	113.38
1	A	173	CYS	N-CA-C	5.91	118.17	110.43
1	A	168	ASN	N-CA-CB	5.89	119.16	110.56
1	A	199	LEU	N-CA-C	-5.85	106.27	113.41
1	A	6	VAL	N-CA-C	5.75	116.08	107.51
1	A	176	THR	CB-CA-C	5.67	119.00	109.48
1	A	104	PHE	N-CA-C	5.66	118.44	109.96
1	A	146	GLU	N-CA-C	5.64	118.16	111.33
1	A	60	ASP	N-CA-C	-5.60	102.99	110.55
1	A	312	PRO	CA-C-N	5.59	132.22	121.54
1	A	312	PRO	C-N-CA	5.59	132.22	121.54
1	A	282	LYS	N-CA-C	5.59	118.45	111.24
1	A	221	GLU	N-CA-C	5.56	122.64	110.80
1	A	8	TRP	CA-C-N	-5.52	114.30	122.41
1	A	8	TRP	C-N-CA	-5.52	114.30	122.41
1	A	90	THR	CB-CA-C	-5.51	103.48	111.14
1	A	84	THR	N-CA-C	-5.46	101.06	109.52
1	A	150	ALA	N-CA-C	-5.43	105.44	111.36
1	A	313	ARG	NE-CZ-NH2	5.40	124.06	119.20
1	A	252	SER	N-CA-C	5.38	117.47	110.43
1	A	160	GLY	N-CA-C	-5.37	108.09	114.48
1	A	86	ARG	N-CA-C	-5.36	105.52	111.36
1	A	6	VAL	CB-CA-C	-5.09	104.18	111.31
1	A	105	GLN	N-CA-C	-5.05	102.71	110.14
1	A	256	VAL	CB-CA-C	5.03	118.14	110.90
1	A	309	ARG	O-C-N	5.03	128.99	122.81

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	70	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	70	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 (70) 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:119:LEU:HG	1:A:120:ARG:HG3	1.80	0.63
1:A:26:MET:HE3	1:A:32:PRO:HD2	1.80	0.62
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
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:7:GLN:HA	1:A:35:SER:HB3	1.94	0.50
1:A:51:GLU:HG3	1:A:53:ARG:HG3	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:97:VAL:HG21	1:A:209:ILE:HD11	1.96	0.47
1:A:313:ARG:HA	1:A:313:ARG:HD2	1.75	0.47
1:A:32:PRO:HD3	1:A:273:GLN:CD	2.39	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:HH11	1:A:224:ARG:HD3	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:107:ASN:HD22	1:A:107:ASN:H	1.65	0.44
1:A:278:PHE:CD1	1:A:286:PHE:HB3	2.52	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
1:A:118:GLY:HA3	1:A:121:GLU:HG3	2.00	0.42
1:A:269:ASN:OD1	1:A:272:ARG:NH1	2.53	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:49:ILE:HD11	1:A:56:ALA:C	2.46	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:OD1	1:A:240:ASP:N	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%)	9 18

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%)	204 (77%)	60 (23%)	1 1

All (60) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	SER
1	A	23	GLN
1	A	24	ARG
1	A	26	MET
1	A	29	VAL
1	A	39	ARG
1	A	43	ILE
1	A	49	ILE
1	A	52	ASN
1	A	75	ARG
1	A	85	GLU
1	A	87	GLN
1	A	89	ARG
1	A	97	VAL
1	A	99	LYS
1	A	105	GLN
1	A	106	LEU
1	A	107	ASN
1	A	109	LEU
1	A	113	LYS
1	A	121	GLU
1	A	132	LEU
1	A	136	LEU
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	186	GLN
1	A	187	GLU
1	A	191	SER
1	A	197	LYS
1	A	211	GLU
1	A	214	VAL

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Mol	Chain	Res	Type
1	A	218	LEU
1	A	219	SER
1	A	220	ASP
1	A	224	ARG
1	A	226	GLU
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

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	13	GLN
1	A	23	GLN
1	A	87	GLN
1	A	91	HIS
1	A	107	ASN
1	A	126	ASN
1	A	179	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 [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 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	3,3,3	0.42	0	2,3,3	0.94	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 [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](#)

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%)	-0.57	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 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(Å ²)	Q<0.9
3	CO3	A	401	4/4	0.95	0.09	22,26,31,33	0
2	FE	A	400	1/1	1.00	0.01	23,23,23,23	0

6.5 Other polymers [i](#)

There are no such residues in this entry.