



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 11:26 AM UTC

PDB ID : 1VLZ / pdb_00001vlz
Title : UNCOUPLED PHOSPHORYLATION AND ACTIVATION IN BACTERIAL
CHEMOTAXIS: THE 2.1 ANGSTROM STRUCTURE OF A THREONINE
TO ISOLEUCINE MUTANT AT POSITION 87 OF CHEY
Authors : Ganguli, S.; Wang, H.; Matsumura, P.; Volz, K.
Deposited on : 1995-04-21
Resolution : 2.05 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
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

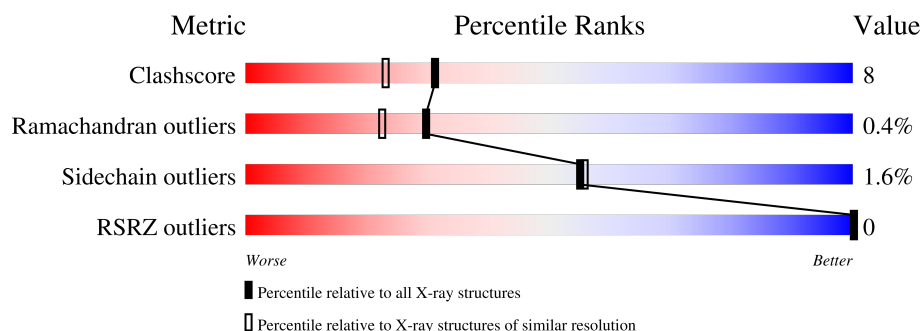
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.05 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	2333 (2.04-2.04)
Ramachandran outliers	187476	2318 (2.04-2.04)
Sidechain outliers	187428	2318 (2.04-2.04)
RSRZ outliers	180081	2260 (2.04-2.04)

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	128	60% 33% 5% .
1	B	128	52% 38% 9% .

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 2070 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 CHEY.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	128	Total	C	N	O	S	0	0	0
			954	609	156	183	6			
1	B	127	Total	C	N	O	S	0	0	0
			944	604	158	176	6			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	87	ILE	THR	conflict	UNP P06143
B	87	ILE	THR	conflict	UNP P06143

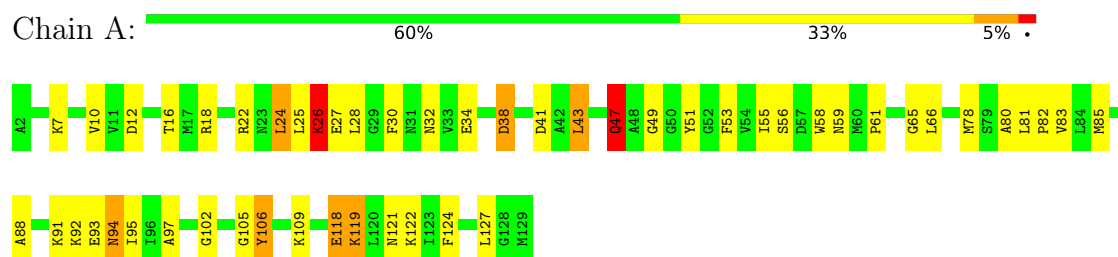
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	92	Total	O	0	0
			92	92		
2	B	80	Total	O	0	0
			80	80		

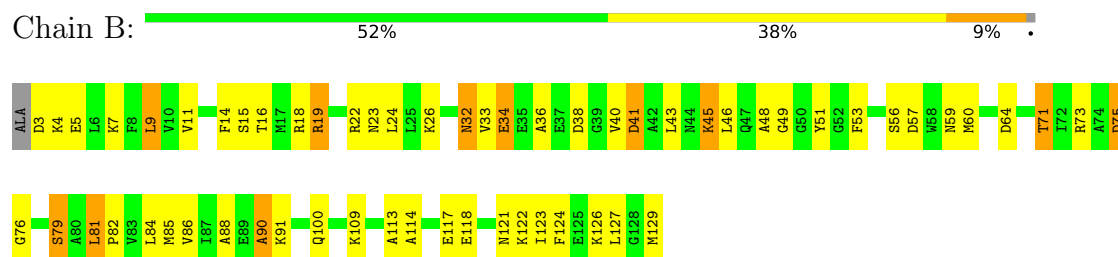
3 Residue-property plots

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: CHEY



• Molecule 1: CHEY



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	54.04Å 72.47Å 36.15Å 90.00° 109.06° 90.00°	Depositor
Resolution (Å)	10.00 – 2.05 10.00 – 2.05	Depositor EDS
% Data completeness (in resolution range)	(Not available) (10.00-2.05) 86.9 (10.00-2.05)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.37 (at 2.05Å)	Xtriage
Refinement program	PROFFT	Depositor
R, R_{free}	0.156 , (Not available) 0.162 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	17.6	Xtriage
Anisotropy	0.073	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 89.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.048 for -h-l,-k,l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	2070	wwPDB-VP
Average B, all atoms (Å ²)	17.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 10.95% 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

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	1.14	0/966	2.42	54/1304 (4.1%)
1	B	1.14	0/956	2.37	69/1290 (5.3%)
All	All	1.14	0/1922	2.39	123/2594 (4.7%)

There are no bond length outliers.

All (123) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	18	ARG	NE-CZ-NH1	12.07	133.57	121.50
1	A	47	GLN	OE1-CD-NE2	10.80	133.40	122.60
1	A	94	ASN	OD1-CG-ND2	10.35	132.95	122.60
1	B	49	GLY	N-CA-C	9.27	127.44	112.61
1	B	100	GLN	OE1-CD-NE2	-9.02	113.58	122.60
1	B	22	ARG	NE-CZ-NH1	-8.65	112.85	121.50
1	B	32	ASN	N-CA-CB	-8.40	96.58	110.53
1	B	126	LYS	CA-CB-CG	8.32	130.74	114.10
1	B	19	ARG	CD-NE-CZ	8.24	135.94	124.40
1	A	47	GLN	CA-C-O	-7.94	111.13	120.10
1	B	60	MET	O-C-N	7.92	127.40	121.85
1	A	18	ARG	NE-CZ-NH2	-7.88	112.11	119.20
1	B	59	ASN	CA-CB-CG	7.75	120.35	112.60
1	A	119	LYS	CA-C-O	-7.59	112.37	120.42
1	A	91	LYS	CB-CA-C	-7.43	99.37	110.67
1	A	26	LYS	CA-C-N	7.43	131.60	120.31
1	A	26	LYS	C-N-CA	7.43	131.60	120.31
1	A	32	ASN	OD1-CG-ND2	7.39	129.99	122.60
1	A	95	ILE	N-CA-C	-7.34	103.63	110.53
1	A	24	LEU	CA-C-N	7.29	130.05	120.28
1	A	24	LEU	C-N-CA	7.29	130.05	120.28
1	B	76	GLY	CA-C-N	7.17	130.38	120.63
1	B	76	GLY	C-N-CA	7.17	130.38	120.63
1	B	32	ASN	CB-CA-C	7.10	123.54	111.68
1	B	23	ASN	CA-CB-CG	-7.07	105.53	112.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	100	GLN	CA-CB-CG	7.04	128.18	114.10
1	B	45	LYS	CA-CB-CG	-6.97	100.16	114.10
1	B	81	LEU	N-CA-CB	-6.92	97.35	109.73
1	A	91	LYS	CA-C-O	-6.88	112.44	120.62
1	A	102	GLY	CA-C-O	-6.84	110.85	119.19
1	B	73	ARG	NE-CZ-NH2	-6.78	113.10	119.20
1	B	23	ASN	CA-C-O	-6.69	113.32	120.42
1	A	24	LEU	CA-C-O	-6.68	113.81	120.82
1	B	22	ARG	NE-CZ-NH2	6.67	125.21	119.20
1	A	119	LYS	N-CA-CB	6.65	120.00	110.16
1	A	97	ALA	CA-C-N	6.64	129.07	120.44
1	A	97	ALA	C-N-CA	6.64	129.07	120.44
1	A	38	ASP	CA-C-O	-6.63	114.19	121.28
1	B	26	LYS	N-CA-CB	6.57	119.72	109.94
1	A	93	GLU	CB-CG-CD	6.49	123.63	112.60
1	B	41	ASP	CB-CA-C	6.45	121.50	110.79
1	A	118	GLU	CG-CD-OE2	-6.43	103.62	118.40
1	A	88	ALA	N-CA-C	6.42	120.85	112.89
1	B	53	PHE	CA-CB-CG	6.40	120.20	113.80
1	B	109	LYS	N-CA-CB	6.40	117.77	110.03
1	B	56	SER	N-CA-CB	-6.26	99.66	111.00
1	A	53	PHE	CA-CB-CG	6.25	120.05	113.80
1	A	10	VAL	CA-C-O	-6.15	113.95	120.53
1	A	80	ALA	CA-C-O	-6.13	112.39	119.56
1	B	14	PHE	CA-C-O	-6.11	114.24	120.84
1	B	49	GLY	CA-C-O	6.02	128.85	122.59
1	A	49	GLY	CA-C-O	6.02	127.28	122.52
1	B	9	LEU	CA-C-N	6.02	130.61	122.90
1	B	9	LEU	C-N-CA	6.02	130.61	122.90
1	A	106	TYR	CA-CB-CG	-6.00	103.10	113.90
1	A	78	MET	CA-CB-CG	5.99	126.08	114.10
1	B	24	LEU	CA-C-N	5.96	128.19	120.44
1	B	24	LEU	C-N-CA	5.96	128.19	120.44
1	A	56	SER	CA-C-O	-5.95	113.81	120.66
1	A	16	THR	CA-CB-OG1	-5.95	100.68	109.60
1	B	71	THR	CA-CB-OG1	-5.89	100.76	109.60
1	B	75	ASP	CB-CG-OD1	5.87	131.90	118.40
1	B	57	ASP	CA-CB-CG	-5.87	106.73	112.60
1	A	22	ARG	CD-NE-CZ	-5.84	116.22	124.40
1	A	66	LEU	CB-CA-C	5.83	120.47	110.79
1	A	56	SER	N-CA-C	5.83	118.97	109.24
1	B	64	ASP	CA-C-N	5.83	126.45	119.98

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	64	ASP	C-N-CA	5.83	126.45	119.98
1	B	79	SER	N-CA-C	5.80	118.35	111.33
1	B	40	VAL	O-C-N	5.79	127.93	121.90
1	B	48	ALA	CA-C-N	5.79	129.03	121.85
1	B	48	ALA	C-N-CA	5.79	129.03	121.85
1	B	16	THR	CA-C-O	-5.77	114.43	120.55
1	A	43	LEU	CA-C-O	-5.74	114.47	120.55
1	A	121	ASN	CA-C-O	-5.73	114.47	120.55
1	A	55	ILE	CA-C-N	-5.72	113.46	122.73
1	A	55	ILE	C-N-CA	-5.72	113.46	122.73
1	B	84	LEU	CA-C-O	-5.67	114.08	120.32
1	A	30	PHE	O-C-N	5.66	129.79	123.05
1	A	7	LYS	CG-CD-CE	5.64	124.27	111.30
1	B	11	VAL	O-C-N	5.58	129.12	123.20
1	A	41	ASP	CB-CG-OD2	5.56	131.20	118.40
1	B	109	LYS	N-CA-C	-5.50	102.71	110.29
1	A	59	ASN	N-CA-C	5.49	117.70	107.99
1	A	65	GLY	CA-C-N	5.48	127.62	120.28
1	A	65	GLY	C-N-CA	5.48	127.62	120.28
1	A	122	LYS	N-CA-CB	5.45	118.14	110.12
1	B	19	ARG	NE-CZ-NH2	-5.45	114.30	119.20
1	B	19	ARG	NE-CZ-NH1	5.43	126.93	121.50
1	B	34	GLU	N-CA-CB	5.43	120.00	111.20
1	B	75	ASP	CA-CB-CG	5.43	118.03	112.60
1	B	113	ALA	N-CA-C	-5.41	105.49	111.71
1	B	16	THR	O-C-N	5.41	127.85	122.12
1	B	117	GLU	CG-CD-OE2	-5.33	106.13	118.40
1	B	38	ASP	O-C-N	5.33	128.69	122.99
1	A	91	LYS	O-C-N	5.33	129.96	122.77
1	A	92	LYS	CB-CA-C	-5.32	101.90	110.74
1	A	83	VAL	CA-CB-CG2	5.31	119.43	110.40
1	B	4	LYS	N-CA-C	-5.29	106.89	113.19
1	B	114	ALA	O-C-N	5.29	127.52	122.07
1	B	88	ALA	N-CA-C	5.28	119.44	112.89
1	A	109	LYS	N-CA-CB	5.28	116.41	110.03
1	B	123	ILE	N-CA-C	-5.25	105.59	110.53
1	B	18	ARG	N-CA-CB	5.24	117.91	110.16
1	B	121	ASN	CB-CA-C	5.22	119.73	110.85
1	B	46	LEU	O-C-N	5.18	127.61	122.12
1	B	113	ALA	O-C-N	5.14	128.23	122.22
1	B	45	LYS	CA-C-N	5.13	127.16	120.28
1	B	45	LYS	C-N-CA	5.13	127.16	120.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	100	GLN	CG-CD-NE2	5.11	124.07	116.40
1	B	33	VAL	N-CA-CB	-5.10	104.07	111.52
1	B	91	LYS	N-CA-CB	5.10	117.56	109.71
1	A	124	PHE	CA-C-O	-5.09	115.47	121.07
1	A	105	GLY	CA-C-O	-5.09	116.63	121.62
1	B	86	VAL	CA-C-O	-5.08	115.05	120.59
1	B	45	LYS	N-CA-C	5.06	117.91	111.69
1	B	90	ALA	CA-C-O	-5.05	113.28	120.51
1	A	119	LYS	O-C-N	5.04	127.89	122.15
1	B	126	LYS	N-CA-CB	5.04	118.09	110.28
1	A	12	ASP	CB-CG-OD2	5.04	129.99	118.40
1	B	45	LYS	N-CA-CB	-5.03	102.71	110.20
1	A	27	GLU	CG-CD-OE2	-5.02	106.86	118.40
1	B	38	ASP	N-CA-CB	5.01	119.15	111.24

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	954	0	951	16	0
1	B	944	0	947	15	0
2	A	92	0	0	3	0
2	B	80	0	0	0	0
All	All	2070	0	1898	31	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 8.

All (31) 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:B:127:LEU:HD23	1:B:129:MET:SD	2.12	0.89
1:A:58:TRP:CE3	1:A:85:MET:HE2	2.21	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:75:ASP:O	1:B:79:SER:HB3	1.95	0.66
1:A:34:GLU:OE1	2:A:151:HOH:O	2.14	0.65
1:A:58:TRP:HE3	1:A:85:MET:HE2	1.68	0.57
1:A:43:LEU:O	1:A:47:GLN:HG2	2.04	0.56
1:B:43:LEU:HD11	1:B:71:THR:HG21	1.88	0.56
1:B:118:GLU:O	1:B:122:LYS:HG3	2.06	0.56
1:A:94:ASN:ND2	2:A:196:HOH:O	2.16	0.55
1:B:3:ASP:C	1:B:5:GLU:H	2.15	0.54
1:B:34:GLU:HG3	1:B:51:TYR:OH	2.08	0.53
1:B:127:LEU:HB3	1:B:129:MET:HG3	1.91	0.53
1:B:15:SER:HB2	1:B:19:ARG:HH11	1.76	0.49
1:A:106:TYR:C	1:A:106:TYR:CD1	2.92	0.47
1:A:38:ASP:HB2	1:A:61:PRO:O	2.15	0.46
1:B:85:MET:HE3	1:B:85:MET:HB3	1.63	0.45
1:A:118:GLU:HG2	2:A:212:HOH:O	2.16	0.45
1:B:124:PHE:HD1	1:B:129:MET:SD	2.40	0.45
1:B:3:ASP:C	1:B:5:GLU:N	2.76	0.43
1:A:26:LYS:HB2	1:A:26:LYS:HE3	0.97	0.43
1:A:34:GLU:HG3	1:A:51:TYR:OH	2.19	0.43
1:B:81:LEU:HA	1:B:82:PRO:HD3	1.88	0.43
1:A:24:LEU:O	1:A:28:LEU:HG	2.19	0.43
1:B:41:ASP:O	1:B:45:LYS:HG3	2.18	0.42
1:A:85:MET:O	1:A:106:TYR:HA	2.18	0.42
1:A:81:LEU:HA	1:A:82:PRO:HD3	1.95	0.42
1:A:106:TYR:O	1:A:119:LYS:HE2	2.19	0.42
1:B:9:LEU:HD11	1:B:36:ALA:HB2	2.03	0.41
1:B:7:LYS:HE2	1:B:32:ASN:OD1	2.20	0.41
1:A:25:LEU:HA	1:A:25:LEU:HD23	1.81	0.40
1:A:47:GLN:HE21	1:A:47:GLN:HA	1.87	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	126/128 (98%)	124 (98%)	2 (2%)	0	100	100
1	B	125/128 (98%)	121 (97%)	3 (2%)	1 (1%)	16	9
All	All	251/256 (98%)	245 (98%)	5 (2%)	1 (0%)	30	22

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	90	ALA

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	96/102 (94%)	93 (97%)	3 (3%)	35	30
1	B	94/102 (92%)	94 (100%)	0	100	100
All	All	190/204 (93%)	187 (98%)	3 (2%)	55	56

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	26	LYS
1	A	47	GLN
1	A	127	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	47	GLN
1	A	59	ASN

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](#)

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](#)

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	128/128 (100%)	-0.79	0 100 100	7, 12, 25, 34	0
1	B	127/128 (99%)	-0.61	0 100 100	7, 17, 31, 50	0
All	All	255/256 (99%)	-0.70	0 100 100	7, 14, 30, 50	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.