



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 10:30 PM UTC

PDB ID : 3ZKW / pdb_00003zkw
Title : Periplasmic Binding Protein CeuE apo form
Authors : Raines, D.J.; Moroz, O.V.; Wilson, K.S.; Duhme-Klair, A.K.
Deposited on : 2013-01-25
Resolution : 1.71 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

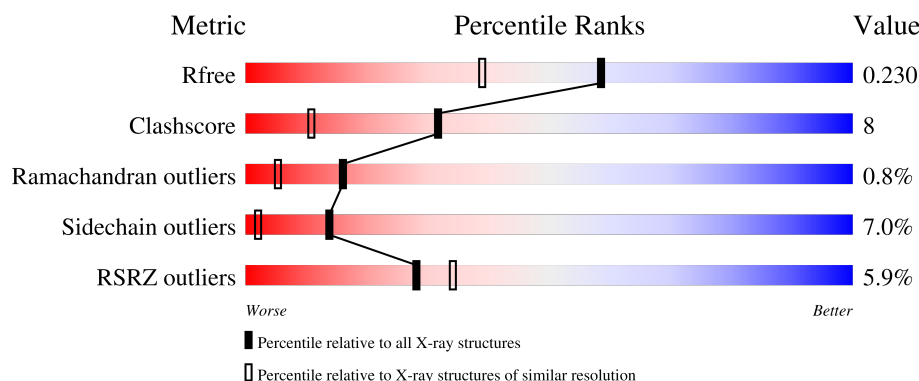
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.71 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))
R_{free}	180053	1039 (1.72-1.72)
Clashscore	190562	1049 (1.72-1.72)
Ramachandran outliers	187476	1041 (1.72-1.72)
Sidechain outliers	187428	1041 (1.72-1.72)
RSRZ outliers	180081	1039 (1.72-1.72)

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	287	8% 83% 12% ..
1	C	287	6% 80% 14% ..
2	B	287	5% 81% 16% ..

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6763 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 ENTEROCHELIN UPTAKE PERIPLASMIC BINDING PROTEIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	287	Total	C	N	O	S	48	0	0
			2220	1428	363	427	2			
1	C	286	Total	C	N	O	S	17	0	0
			2215	1424	363	426	2			

- Molecule 2 is a protein called ENTEROCHELIN UPTAKE PERIPLASMIC BINDING PROTEIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	287	Total	C	N	O	S	19	0	0
			2220	1428	363	427	2			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	24	ILE	LEU	conflict	UNP Q0P8Q4
B	41	ILE	LEU	conflict	UNP Q0P8Q4

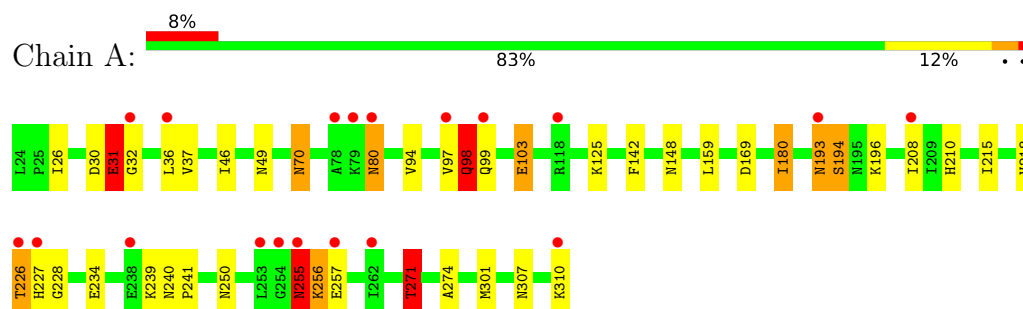
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	41	Total	O	0	0
			41	41		
3	B	38	Total	O	0	0
			38	38		
3	C	29	Total	O	0	0
			29	29		

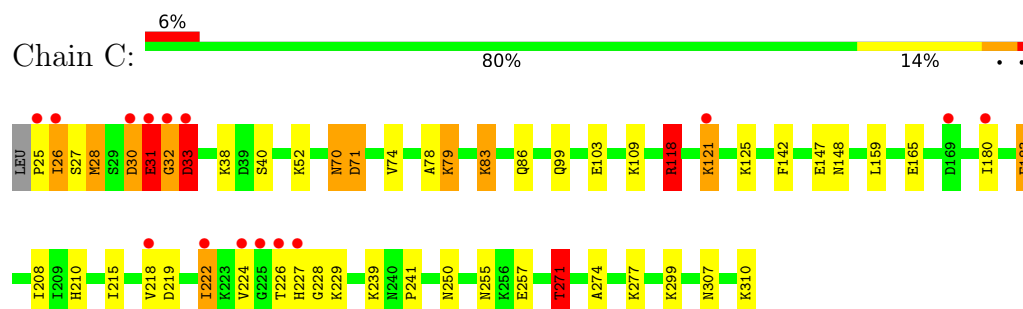
3 Residue-property plots [i](#)

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.

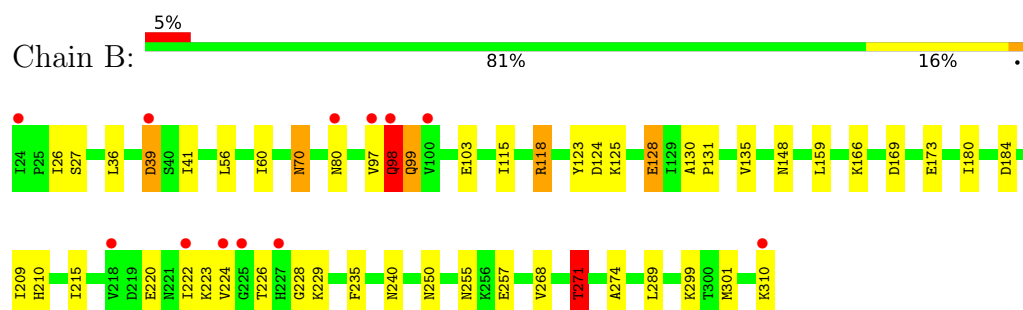
• Molecule 1: ENTEROCHELIN UPTAKE PERIPLASMIC BINDING PROTEIN



• Molecule 1: ENTEROCHELIN UPTAKE PERIPLASMIC BINDING PROTEIN



• Molecule 2: ENTEROCHELIN UPTAKE PERIPLASMIC BINDING PROTEIN



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	56.95Å 62.74Å 67.98Å 82.19° 76.74° 75.96°	Depositor
Resolution (Å)	65.93 – 1.71 65.93 – 1.71	Depositor EDS
% Data completeness (in resolution range)	97.0 (65.93-1.71) 97.0 (65.93-1.71)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.13 (at 1.71Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.195 , 0.227 0.198 , 0.230	Depositor DCC
R_{free} test set	4647 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	25.6	Xtriage
Anisotropy	0.046	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 24.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	6763	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.81% 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	2.38	16/2253 (0.7%)	1.54	26/3044 (0.9%)
1	C	1.42	14/2248 (0.6%)	1.22	16/3035 (0.5%)
2	B	1.55	8/2253 (0.4%)	1.36	11/3044 (0.4%)
All	All	1.83	38/6754 (0.6%)	1.38	53/9123 (0.6%)

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	2
1	C	0	1
2	B	1	2
All	All	1	5

All (38) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	97	VAL	CA-C	73.13	2.33	1.52
1	A	97	VAL	C-O	-36.33	0.88	1.24
2	B	118	ARG	NE-CZ	28.62	1.64	1.33
1	A	80	ASN	CA-CB	28.16	2.02	1.53
2	B	99	GLN	CA-CB	-24.11	1.15	1.53
2	B	184	ASP	CA-CB	-21.50	1.18	1.53
1	A	256	LYS	C-N	-20.13	1.06	1.33
1	A	103	GLU	CG-CD	-19.80	1.02	1.52
1	A	97	VAL	CA-CB	19.64	1.77	1.53
1	C	147	GLU	CD-OE1	-17.59	0.92	1.25
1	C	147	GLU	CD-OE2	16.52	1.56	1.25
1	C	31	GLU	N-CA	15.96	1.66	1.46
1	A	227	HIS	CB-CG	-15.26	1.28	1.50
1	A	234	GLU	CG-CD	-14.38	1.16	1.52

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	97	VAL	C-N	12.46	1.51	1.33
1	C	226	THR	CA-CB	-10.88	1.35	1.53
2	B	98	GLN	CA-CB	10.23	1.70	1.53
1	C	31	GLU	CA-C	8.32	1.64	1.52
1	A	255	ASN	C-N	-8.30	1.21	1.33
1	C	30	ASP	CA-C	7.66	1.63	1.52
1	C	33	ASP	CA-C	6.80	1.61	1.52
2	B	240	ASN	N-CA	-6.70	1.41	1.46
1	C	299	LYS	CG-CD	6.68	1.72	1.52
1	C	83	LYS	CG-CD	6.57	1.72	1.52
1	A	49	ASN	CA-CB	6.51	1.57	1.52
1	A	240	ASN	CA-C	-6.41	1.48	1.53
1	C	30	ASP	N-CA	6.36	1.54	1.46
1	C	109	LYS	C-O	-6.18	1.21	1.23
1	A	226	THR	CA-CB	-6.06	1.43	1.53
1	A	98	GLN	CA-CB	-5.95	1.43	1.53
1	C	33	ASP	N-CA	5.88	1.53	1.46
1	A	196	LYS	CG-CD	5.62	1.69	1.52
1	C	165	GLU	CA-CB	5.39	1.61	1.53
2	B	235	PHE	C-O	-5.36	1.17	1.24
2	B	39	ASP	CB-CG	5.34	1.65	1.52
2	B	240	ASN	C-O	-5.32	1.21	1.23
1	C	71	ASP	CG-OD1	5.12	1.35	1.25
1	A	301	MET	SD-CE	-5.01	1.67	1.79

All (53) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	97	VAL	CB-CA-C	-27.96	73.87	111.33
1	A	103	GLU	CB-CG-CD	23.52	152.59	112.60
2	B	118	ARG	CD-NE-CZ	-23.09	92.07	124.40
1	A	97	VAL	N-CA-C	-21.57	80.36	108.84
2	B	118	ARG	NE-CZ-NH1	-15.93	105.57	121.50
1	A	255	ASN	O-C-N	-15.64	101.78	122.59
2	B	99	GLN	CB-CA-C	15.64	134.24	110.62
2	B	98	GLN	CA-CB-CG	12.59	139.28	114.10
2	B	118	ARG	NE-CZ-NH2	12.30	130.27	119.20
1	A	226	THR	N-CA-CB	11.69	128.40	110.28
1	A	80	ASN	N-CA-CB	-11.67	91.52	110.42
1	C	83	LYS	CG-CD-CE	11.51	137.77	111.30
1	A	31	GLU	N-CA-C	10.92	125.67	110.06
1	A	193	ASN	N-CA-C	10.63	129.56	113.61

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	234	GLU	CB-CG-CD	10.42	130.31	112.60
1	A	301	MET	CG-SD-CE	-9.80	79.33	100.90
1	C	226	THR	N-CA-CB	-8.92	96.95	110.16
2	B	99	GLN	CA-CB-CG	-8.83	96.44	114.10
1	A	97	VAL	CA-CB-CG2	-8.79	95.45	110.40
1	A	196	LYS	CB-CG-CD	-8.72	91.25	111.30
2	B	271	THR	OG1-CB-CG2	8.43	126.17	109.30
1	A	97	VAL	N-CA-CB	-8.30	101.61	110.82
1	A	98	GLN	N-CA-CB	8.12	124.22	110.49
1	C	30	ASP	N-CA-C	7.95	127.73	110.80
2	B	301	MET	CG-SD-CE	-7.77	83.81	100.90
1	A	97	VAL	CA-C-N	-7.74	106.76	121.54
1	A	97	VAL	C-N-CA	-7.74	106.76	121.54
1	C	147	GLU	CG-CD-OE2	-7.67	100.76	118.40
1	C	227	HIS	N-CA-C	7.29	122.30	113.41
1	A	271	THR	N-CA-CB	-7.26	99.72	110.17
2	B	271	THR	N-CA-CB	-7.13	99.46	110.29
1	C	271	THR	N-CA-CB	-6.80	99.95	110.29
1	C	31	GLU	N-CA-C	6.79	125.26	110.80
2	B	98	GLN	CB-CA-C	6.24	122.85	110.42
1	A	97	VAL	O-C-N	6.21	129.80	122.66
1	C	165	GLU	CA-CB-CG	6.04	126.18	114.10
1	A	196	LYS	CG-CD-CE	-5.95	97.61	111.30
1	A	226	THR	CA-CB-OG1	-5.90	100.75	109.60
1	C	147	GLU	CG-CD-OE1	5.87	131.91	118.40
1	A	226	THR	CA-CB-CG2	5.80	120.36	110.50
1	A	97	VAL	CA-CB-CG1	-5.79	100.56	110.40
1	A	94	VAL	N-CA-C	-5.63	107.25	112.43
1	C	226	THR	N-CA-C	5.60	117.47	111.36
2	B	184	ASP	CB-CA-C	5.59	120.94	110.70
1	A	80	ASN	CA-CB-CG	-5.57	107.03	112.60
1	C	118	ARG	CB-CG-CD	5.54	124.04	111.30
1	A	271	THR	OG1-CB-CG2	5.33	119.97	109.30
1	A	169	ASP	CB-CA-C	-5.33	102.51	110.88
1	C	31	GLU	O-C-N	-5.25	115.61	122.59
1	C	299	LYS	CB-CG-CD	-5.22	99.29	111.30
1	C	118	ARG	CG-CD-NE	5.21	123.45	112.00
1	C	31	GLU	CA-C-N	5.17	131.01	121.70
1	C	31	GLU	C-N-CA	5.17	131.01	121.70

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	B	98	GLN	CA

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	255	ASN	Mainchain
1	A	80	ASN	Peptide
2	B	118	ARG	Sidechain
2	B	223	LYS	Peptide
1	C	32	GLY	Peptide

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	2220	0	2272	26	0
1	C	2215	0	2275	46	0
2	B	2220	0	2276	38	0
3	A	41	0	0	0	0
3	B	38	0	0	3	0
3	C	29	0	0	2	0
All	All	6763	0	6823	108	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 (108) 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:C:224:VAL:HG21	1:C:228:GLY:HA2	1.44	0.99
1:A:26:ILE:H	1:A:148:ASN:HD21	1.07	0.96
1:A:250:ASN:HD22	1:A:255:ASN:HD22	1.18	0.91
1:A:30:ASP:O	1:A:31:GLU:OE1	1.90	0.89
2:B:26:ILE:H	2:B:148:ASN:HD21	1.19	0.89
1:C:224:VAL:HG21	1:C:228:GLY:CA	2.03	0.88
1:C:78:ALA:O	1:C:79:LYS:CB	2.20	0.86
1:C:27:SER:H	1:C:148:ASN:HD21	1.26	0.84

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:27:SER:HB2	3:C:2001:HOH:O	1.82	0.77
1:A:142:PHE:CE2	1:A:208:ILE:HD11	2.20	0.76
1:C:250:ASN:HD22	1:C:255:ASN:HD22	1.33	0.76
1:C:218:VAL:HG11	1:C:241:PRO:HB3	1.69	0.75
1:A:250:ASN:ND2	1:A:255:ASN:HD22	1.85	0.74
1:A:224:VAL:HG13	1:A:225:GLY:O	1.89	0.72
1:A:193:ASN:O	1:A:194:SER:HB3	1.89	0.71
2:B:250:ASN:HD22	2:B:255:ASN:HD22	1.39	0.71
1:C:78:ALA:O	1:C:79:LYS:HB3	1.91	0.70
1:A:250:ASN:HD22	1:A:255:ASN:ND2	1.90	0.69
1:C:25:PRO:HB2	1:C:40:SER:HB3	1.74	0.69
2:B:27:SER:HB2	1:C:121:LYS:HG3	1.74	0.69
2:B:70:ASN:H	2:B:70:ASN:HD22	1.43	0.66
1:C:28:MET:HE2	1:C:38:LYS:HG3	1.75	0.66
1:C:224:VAL:CG2	1:C:228:GLY:H	2.08	0.66
1:C:27:SER:CB	3:C:2001:HOH:O	2.44	0.63
1:C:28:MET:HE1	1:C:30:ASP:OD2	1.98	0.62
2:B:97:VAL:HG11	2:B:289:LEU:HD22	1.82	0.62
1:C:224:VAL:HG21	1:C:228:GLY:N	2.14	0.62
1:C:83:LYS:HA	1:C:86:GLN:HG2	1.80	0.62
1:C:194:SER:H	1:C:250:ASN:HD21	1.49	0.60
1:A:70:ASN:H	1:A:70:ASN:HD22	1.49	0.60
1:A:222:ILE:O	1:A:223:LYS:O	2.19	0.60
1:C:219:ASP:CG	1:C:222:ILE:HD13	2.26	0.60
2:B:26:ILE:N	2:B:148:ASN:HD21	1.94	0.59
2:B:26:ILE:H	2:B:148:ASN:ND2	1.94	0.58
2:B:80:ASN:OD1	2:B:97:VAL:HG22	2.03	0.58
1:A:26:ILE:N	1:A:148:ASN:HD21	1.91	0.58
1:C:250:ASN:ND2	1:C:255:ASN:HD22	2.00	0.58
1:A:37:VAL:HG21	1:A:46:ILE:HD12	1.85	0.57
2:B:41:ILE:HG22	2:B:41:ILE:O	2.05	0.56
2:B:166:LYS:O	2:B:169:ASP:OD1	2.24	0.56
2:B:41:ILE:HG21	2:B:123:TYR:CG	2.41	0.56
1:A:30:ASP:C	1:A:31:GLU:OE1	2.49	0.55
1:C:70:ASN:HD22	1:C:70:ASN:H	1.54	0.55
1:C:307:ASN:ND2	1:C:310:LYS:NZ	2.54	0.55
2:B:173:GLU:HG3	2:B:299:LYS:HE2	1.88	0.55
1:A:222:ILE:O	1:A:223:LYS:C	2.49	0.55
2:B:194:SER:H	2:B:250:ASN:HD21	1.54	0.54
1:A:26:ILE:H	1:A:148:ASN:ND2	1.90	0.54
1:C:32:GLY:O	1:C:33:ASP:HB3	2.07	0.54

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:250:ASN:ND2	2:B:255:ASN:HD22	2.05	0.54
2:B:56:LEU:HD12	2:B:115:ILE:HG22	1.91	0.53
1:A:271:THR:CG2	1:A:274:ALA:H	2.21	0.53
1:C:26:ILE:HG22	1:C:148:ASN:CG	2.34	0.53
2:B:41:ILE:HG22	2:B:123:TYR:CE2	2.44	0.53
1:C:78:ALA:O	1:C:79:LYS:HB2	2.07	0.52
1:C:194:SER:N	1:C:250:ASN:HD21	2.07	0.51
2:B:60:ILE:HG13	2:B:135:VAL:HG21	1.92	0.51
2:B:130:ALA:HB1	2:B:131:PRO:HD2	1.92	0.51
2:B:210:HIS:HD2	2:B:215:ILE:O	1.94	0.51
1:C:219:ASP:CG	1:C:222:ILE:CD1	2.83	0.51
2:B:80:ASN:OD1	2:B:97:VAL:N	2.33	0.51
2:B:222:ILE:HD12	2:B:229:LYS:CD	2.42	0.49
1:C:219:ASP:OD2	1:C:222:ILE:HD11	2.12	0.49
2:B:70:ASN:H	2:B:70:ASN:ND2	2.09	0.48
1:C:210:HIS:HD2	1:C:215:ILE:O	1.96	0.48
1:A:180:ILE:C	1:A:180:ILE:HD12	2.39	0.48
1:C:219:ASP:OD2	1:C:222:ILE:CD1	2.62	0.48
1:A:193:ASN:O	1:A:194:SER:CB	2.60	0.47
1:C:222:ILE:HD11	1:C:229:LYS:HZ3	1.78	0.47
1:C:307:ASN:ND2	1:C:310:LYS:HZ3	2.12	0.47
1:C:118:ARG:CG	1:C:118:ARG:HH11	2.28	0.47
1:C:271:THR:CG2	1:C:274:ALA:H	2.27	0.47
1:C:52:LYS:HA	1:C:74:VAL:HG13	1.97	0.47
2:B:222:ILE:HD12	2:B:229:LYS:HD3	1.96	0.47
2:B:194:SER:N	2:B:250:ASN:HD21	2.12	0.47
1:A:271:THR:HG23	1:A:274:ALA:H	1.81	0.46
1:C:26:ILE:HG22	1:C:148:ASN:ND2	2.31	0.46
2:B:268:VAL:O	2:B:271:THR:HB	2.15	0.45
1:C:27:SER:H	1:C:148:ASN:ND2	2.04	0.45
1:C:30:ASP:O	1:C:31:GLU:HB2	2.16	0.45
1:A:210:HIS:HD2	1:A:215:ILE:O	2.00	0.45
2:B:41:ILE:HG22	2:B:123:TYR:CD2	2.52	0.45
2:B:271:THR:CG2	2:B:274:ALA:H	2.30	0.45
1:A:218:VAL:HG11	1:A:241:PRO:HB3	1.99	0.45
1:C:183:GLU:H	1:C:183:GLU:CD	2.25	0.45
2:B:228:GLY:N	3:B:2024:HOH:O	1.72	0.44
1:A:239:LYS:HA	1:A:239:LYS:HD3	1.90	0.44
2:B:124:ASP:O	2:B:128:GLU:OE1	2.36	0.43
1:A:224:VAL:HG11	1:A:228:GLY:HA2	1.99	0.43
1:C:32:GLY:O	1:C:33:ASP:CB	2.65	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:27:SER:CB	1:C:121:LYS:HG3	2.45	0.43
1:C:31:GLU:CD	1:C:32:GLY:HA3	2.44	0.43
2:B:228:GLY:CA	3:B:2024:HOH:O	2.45	0.42
2:B:97:VAL:HG11	2:B:289:LEU:CD2	2.49	0.42
2:B:208:ILE:CG2	2:B:209:ILE:N	2.80	0.42
2:B:41:ILE:CG2	2:B:123:TYR:CD2	3.02	0.42
1:C:142:PHE:CD1	1:C:142:PHE:C	2.97	0.42
1:C:222:ILE:CD1	1:C:222:ILE:N	2.83	0.42
1:A:30:ASP:OD1	1:A:32:GLY:HA2	2.20	0.42
2:B:208:ILE:HD13	2:B:208:ILE:HG21	1.59	0.41
2:B:220:GLU:OE2	3:B:2023:HOH:O	2.22	0.41
2:B:41:ILE:HG21	2:B:41:ILE:HD13	1.85	0.41
2:B:41:ILE:O	2:B:41:ILE:CG2	2.70	0.40
1:C:239:LYS:HA	1:C:239:LYS:HD3	1.97	0.40
1:A:307:ASN:HA	1:A:310:LYS:HG2	2.03	0.40
1:C:219:ASP:HB3	1:C:222:ILE:HD13	2.03	0.40
1:A:30:ASP:C	1:A:30:ASP:OD1	2.64	0.40
1:C:222:ILE:N	1:C:222:ILE:HD12	2.36	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	285/287 (99%)	274 (96%)	8 (3%)	3 (1%)	11	3
1	C	284/287 (99%)	272 (96%)	9 (3%)	3 (1%)	11	3
2	B	285/287 (99%)	276 (97%)	8 (3%)	1 (0%)	30	16
All	All	854/861 (99%)	822 (96%)	25 (3%)	7 (1%)	16	5

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	98	GLN
1	A	223	LYS
1	A	194	SER
2	B	98	GLN
1	C	33	ASP
1	C	79	LYS
1	C	31	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	244/250 (98%)	231 (95%)	13 (5%)	20	3
1	C	244/250 (98%)	223 (91%)	21 (9%)	10	1
2	B	244/250 (98%)	227 (93%)	17 (7%)	14	2
All	All	732/750 (98%)	681 (93%)	51 (7%)	14	2

All (51) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	31	GLU
1	A	36	LEU
1	A	70	ASN
1	A	98	GLN
1	A	99	GLN
1	A	103	GLU
1	A	125	LYS
1	A	159	LEU
1	A	180	ILE
1	A	226	THR
1	A	256	LYS
1	A	257	GLU
1	A	271	THR
2	B	36	LEU
2	B	39	ASP
2	B	70	ASN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
2	B	98	GLN
2	B	99	GLN
2	B	103	GLU
2	B	125	LYS
2	B	128	GLU
2	B	159	LEU
2	B	180	ILE
2	B	191	LEU
2	B	208	ILE
2	B	224	VAL
2	B	226	THR
2	B	257	GLU
2	B	271	THR
2	B	310	LYS
1	C	26	ILE
1	C	28	MET
1	C	31	GLU
1	C	33	ASP
1	C	70	ASN
1	C	71	ASP
1	C	99	GLN
1	C	103	GLU
1	C	118	ARG
1	C	121	LYS
1	C	125	LYS
1	C	159	LEU
1	C	180	ILE
1	C	183	GLU
1	C	191	LEU
1	C	195	ASN
1	C	208	ILE
1	C	222	ILE
1	C	257	GLU
1	C	271	THR
1	C	277	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (26) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	70	ASN
1	A	90	ASN
1	A	99	GLN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	148	ASN
1	A	172	ASN
1	A	210	HIS
1	A	216	ASN
1	A	250	ASN
2	B	70	ASN
2	B	90	ASN
2	B	139	ASN
2	B	148	ASN
2	B	172	ASN
2	B	210	HIS
2	B	216	ASN
2	B	250	ASN
2	B	276	ASN
1	C	70	ASN
1	C	99	GLN
1	C	106	ASN
1	C	148	ASN
1	C	210	HIS
1	C	216	ASN
1	C	250	ASN
1	C	276	ASN
1	C	307	ASN

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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

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](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	A	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A	256:LYS	C	257:GLU	N	1.06

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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	286/287 (99%)	0.29	22 (7%)	19 25	15, 26, 45, 70	14 (4%)
1	C	286/287 (99%)	0.37	16 (5%)	30 36	15, 28, 44, 60	9 (3%)
2	B	287/287 (100%)	0.27	13 (4%)	38 45	17, 26, 42, 64	7 (2%)
All	All	859/861 (99%)	0.31	51 (5%)	28 34	15, 27, 44, 70	30 (3%)

All (51) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	97	VAL	7.3
2	B	310	LYS	6.4
1	A	224	VAL	5.9
2	B	224	VAL	5.1
1	C	222	ILE	5.1
1	A	310	LYS	5.1
1	C	226	THR	4.5
1	A	223	LYS	4.2
1	A	227	HIS	4.2
1	C	224	VAL	3.8
1	A	97	VAL	3.6
1	C	26	ILE	3.4
1	C	32	GLY	3.4
1	C	25	PRO	3.4
1	C	30	ASP	3.3
1	C	31	GLU	3.2
2	B	227	HIS	3.0
2	B	39	ASP	2.9
1	A	208	ILE	2.9
1	A	79	LYS	2.9
1	C	121	LYS	2.8
1	C	227	HIS	2.8
1	A	32	GLY	2.7

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	C	180	ILE	2.7
1	A	254	GLY	2.6
1	C	218	VAL	2.6
1	A	99	GLN	2.5
1	C	225	GLY	2.5
1	A	255	ASN	2.4
1	A	80	ASN	2.4
2	B	218	VAL	2.4
2	B	222	ILE	2.4
1	A	226	THR	2.4
1	C	194	SER	2.3
2	B	225	GLY	2.3
2	B	208	ILE	2.3
1	A	78	ALA	2.2
2	B	80	ASN	2.2
2	B	98	GLN	2.2
1	A	225	GLY	2.2
2	B	24	ILE	2.2
1	C	169	ASP	2.1
2	B	100	VAL	2.1
1	C	33	ASP	2.1
1	A	238	GLU	2.1
1	A	257	GLU	2.1
1	A	36	LEU	2.1
1	A	253	LEU	2.1
1	A	193	ASN	2.0
1	A	118	ARG	2.0
1	A	262	ILE	2.0

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.