



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 6, 2024 – 01:49 pm GMT

PDB ID : 5LWH
Title : CeuE (Y288F variant) a periplasmic protein from *Campylobacter jejuni*.
Authors : Wilde, E.J.; Blagova, E.; Hughes, A.; Raines, D.J.; Moroz, O.V.; Turkenburg, J.; Duhme-Klair, A.-K.; Wilson, K.S.
Deposited on : 2016-09-16
Resolution : 1.47 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 : 4.02b-467
Xtriage (Phenix) : 1.13
EDS : 2.36
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.36

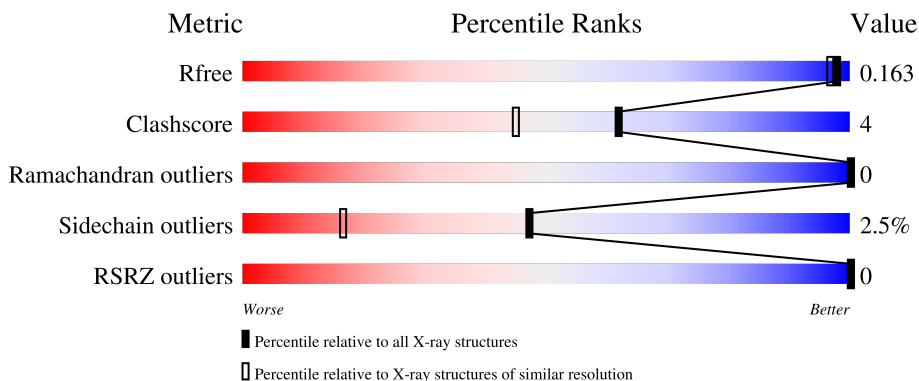
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.47 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}	130704	4690 (1.50-1.46)
Clashscore	141614	4955 (1.50-1.46)
Ramachandran outliers	138981	4846 (1.50-1.46)
Sidechain outliers	138945	4844 (1.50-1.46)
RSRZ outliers	127900	4614 (1.50-1.46)

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	289	81% 16% .

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 2446 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 ABC transporter substrate-binding protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	288	2258	1459	362	434	3	0	5	0

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	22	ALA	-	expression tag	UNP A0A1J6QJN8
A	23	MET	-	expression tag	UNP A0A1J6QJN8
A	288	PHE	TYR	engineered mutation	UNP A0A1J6QJN8

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Zn		
2	A	2	2	2	0	0


- Molecule 3 is water.

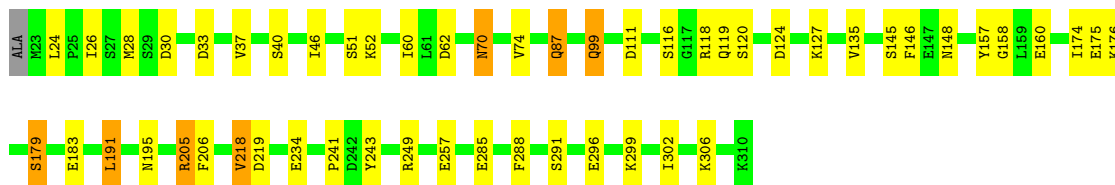
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
3	A	186	186	186	0	0

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 ABC transporter substrate-binding protein

Chain A:  81% 16%



4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	65.52Å 65.52Å 145.65Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	56.81 – 1.47 56.74 – 1.47	Depositor EDS
% Data completeness (in resolution range)	100.0 (56.81-1.47) 100.0 (56.74-1.47)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.99 (at 1.47Å)	Xtrriage
Refinement program	REFMAC 5.8.0158	Depositor
R, R_{free}	0.128 , 0.162 0.129 , 0.163	Depositor DCC
R_{free} test set	3114 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	22.0	Xtrriage
Anisotropy	0.017	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 44.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	0.029 for -h,-k,l	Xtrriage
F_o, F_c correlation	0.98	EDS
Total number of atoms	2446	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.20% 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 i

5.1 Standard geometry i

Bond lengths and bond angles in the following residue types are not validated in this section:
ZN

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	1.49	32/2303 (1.4%)	1.27	19/3110 (0.6%)

All (32) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	160	GLU	CD-OE1	14.11	1.41	1.25
1	A	160	GLU	CG-CD	10.18	1.67	1.51
1	A	120	SER	CB-OG	-9.64	1.29	1.42
1	A	111	ASP	CB-CG	-9.20	1.32	1.51
1	A	40	SER	CB-OG	-8.64	1.31	1.42
1	A	99	GLN	CG-CD	8.39	1.70	1.51
1	A	234	GLU	CD-OE1	7.50	1.33	1.25
1	A	285	GLU	CD-OE1	-7.33	1.17	1.25
1	A	299	LYS	CE-NZ	7.30	1.67	1.49
1	A	51	SER	CB-OG	-7.07	1.33	1.42
1	A	249	ARG	CZ-NH2	-7.00	1.24	1.33
1	A	99	GLN	CD-OE1	6.61	1.38	1.24
1	A	179	SER	CA-CB	6.58	1.62	1.52
1	A	296	GLU	CD-OE2	6.58	1.32	1.25
1	A	160	GLU	CD-OE2	6.43	1.32	1.25
1	A	288	PHE	CG-CD1	-6.38	1.29	1.38
1	A	145	SER	CB-OG	-6.03	1.34	1.42
1	A	52	LYS	CE-NZ	5.98	1.64	1.49
1	A	218	VAL	CB-CG2	-5.93	1.40	1.52
1	A	179	SER	CB-OG	-5.79	1.34	1.42
1	A	257	GLU	CG-CD	5.58	1.60	1.51
1	A	51	SER	CA-CB	-5.56	1.44	1.52
1	A	116	SER	CA-CB	-5.43	1.44	1.52
1	A	145	SER	CA-CB	5.34	1.60	1.52
1	A	257	GLU	CB-CG	-5.29	1.42	1.52
1	A	291	SER	CB-OG	-5.29	1.35	1.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	218	VAL	C-O	5.22	1.33	1.23
1	A	158	GLY	N-CA	5.19	1.53	1.46
1	A	99	GLN	CB-CG	5.18	1.66	1.52
1	A	40	SER	CA-CB	5.13	1.60	1.52
1	A	175	GLU	CD-OE2	-5.09	1.20	1.25
1	A	119	GLN	CG-CD	5.06	1.62	1.51

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	24	LEU	CA-CB-CG	10.23	138.82	115.30
1	A	118	ARG	NE-CZ-NH1	9.90	125.25	120.30
1	A	52	LYS	CD-CE-NZ	7.42	128.78	111.70
1	A	306	LYS	CD-CE-NZ	-7.35	94.79	111.70
1	A	118	ARG	NE-CZ-NH2	-6.95	116.83	120.30
1	A	30	ASP	CB-CG-OD1	-6.91	112.08	118.30
1	A	191[A]	LEU	CB-CG-CD2	6.72	122.42	111.00
1	A	191[B]	LEU	CB-CG-CD2	6.72	122.42	111.00
1	A	28	MET	CG-SD-CE	-6.46	89.87	100.20
1	A	219[A]	ASP	CB-CG-OD2	-6.37	112.56	118.30
1	A	219[B]	ASP	CB-CG-OD2	-6.37	112.56	118.30
1	A	205	ARG	NE-CZ-NH2	-5.84	117.38	120.30
1	A	146	PHE	CB-CG-CD1	-5.71	116.80	120.80
1	A	243	TYR	CG-CD1-CE1	5.61	125.79	121.30
1	A	124	ASP	CB-CG-OD1	5.46	123.21	118.30
1	A	33	ASP	CB-CG-OD1	5.43	123.19	118.30
1	A	179	SER	CA-CB-OG	-5.33	96.80	111.20
1	A	157	TYR	CB-CG-CD2	5.25	124.15	121.00
1	A	249	ARG	NE-CZ-NH2	-5.10	117.75	120.30

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	2258	0	2312	20	0
2	A	2	0	0	0	0
3	A	186	0	0	8	1
All	All	2446	0	2312	20	1

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 4.

All (20) 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:302:ILE:HG12	3:A:670:HOH:O	0.89	1.06
1:A:87:GLN:HE21	1:A:87:GLN:H	1.10	0.92
1:A:205:ARG:HA	3:A:501:HOH:O	1.67	0.92
1:A:26:ILE:H	1:A:148:ASN:HD21	1.24	0.86
1:A:205:ARG:CA	3:A:501:HOH:O	2.21	0.81
1:A:206:PHE:N	3:A:501:HOH:O	1.68	0.81
1:A:302:ILE:CG1	3:A:670:HOH:O	1.68	0.77
1:A:74[A]:VAL:HG21	3:A:577:HOH:O	1.87	0.75
1:A:183:GLU:OE1	3:A:502:HOH:O	2.12	0.66
1:A:70:ASN:HD22	1:A:70:ASN:H	1.43	0.64
1:A:60:ILE:HG13	1:A:135:VAL:HG21	1.81	0.62
1:A:87:GLN:HE21	1:A:87:GLN:N	1.92	0.58
1:A:74[B]:VAL:HG11	3:A:577:HOH:O	2.05	0.55
1:A:174:ILE:HG13	1:A:302:ILE:HD13	1.91	0.51
1:A:218:VAL:HG11	1:A:241:PRO:HB3	1.95	0.48
1:A:176:LYS:O	1:A:179:SER:HB2	2.16	0.46
1:A:37:VAL:HG21	1:A:46:ILE:HD12	2.02	0.42
1:A:70:ASN:HD22	1:A:70:ASN:N	2.13	0.41

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:538:HOH:O	3:A:677:HOH:O[4_545]	2.13	0.07

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	291/289 (101%)	282 (97%)	9 (3%)	0	100 100

There are no Ramachandran outliers to report.

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	246/251 (98%)	240 (98%)	6 (2%)	49 18

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	62	ASP
1	A	70	ASN
1	A	87	GLN
1	A	99	GLN
1	A	127	LYS
1	A	195	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	70	ASN
1	A	87	GLN

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Mol	Chain	Res	Type
1	A	90	ASN
1	A	139	ASN
1	A	148	ASN
1	A	172	ASN
1	A	195	ASN

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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

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	288/289 (99%)	-0.36	0 100 100	17, 25, 42, 59	4 (1%)

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 monosaccharides 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
2	ZN	A	402	1/1	0.98	0.06	38,38,38,38	1
2	ZN	A	401	1/1	1.00	0.09	19,19,19,19	0

6.5 Other polymers [i](#)

There are no such residues in this entry.