wwPDB X-ray Structure Validation Summary Report (i)

Nov 5, 2023 – 10:25 pm GMT

PDB ID	:	8CLZ
Title	:	Crystal structure of Rhizobium etli constitutive L-asparaginase ReAIV (monoclinic form R4mC-2)
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Deposited on	:	2023-02-17
Resolution	:	1.50 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 (i) 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 (i)) were used in the production of this report:

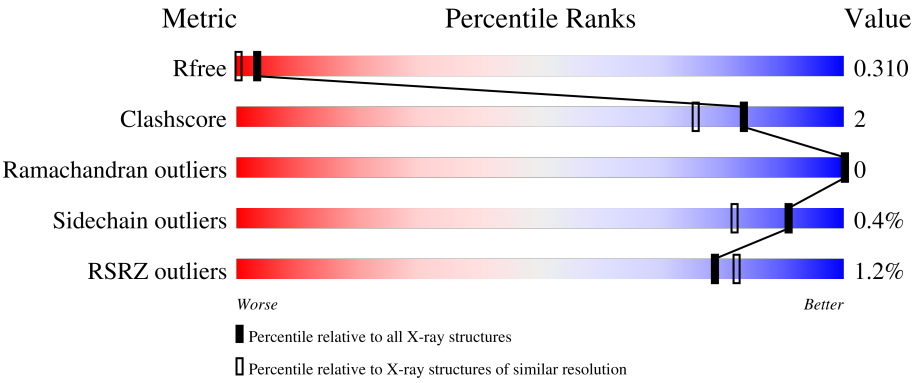
MolProbity	:	4.02b-467
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.50 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))
R_{free}	130704	2936 (1.50-1.50)
Clashscore	141614	3144 (1.50-1.50)
Ramachandran outliers	138981	3066 (1.50-1.50)
Sidechain outliers	138945	3064 (1.50-1.50)
RSRZ outliers	127900	2884 (1.50-1.50)

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 <=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	AAA	341	% 94%
1	BBB	341	% 94%

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 5337 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 Putative L-asparaginase II protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	BBB	334	Total	C	N	O	S	0	3	0
			2482	1544	448	462	28			
1	AAA	333	Total	C	N	O	S	0	3	0
			2471	1538	446	460	27			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
BBB	-5	GLY	-	expression tag	UNP Q2KB35
BBB	-4	ILE	-	expression tag	UNP Q2KB35
BBB	-3	ASP	-	expression tag	UNP Q2KB35
BBB	-2	PRO	-	expression tag	UNP Q2KB35
BBB	-1	PHE	-	expression tag	UNP Q2KB35
BBB	0	THR	-	expression tag	UNP Q2KB35
AAA	-5	GLY	-	expression tag	UNP Q2KB35
AAA	-4	ILE	-	expression tag	UNP Q2KB35
AAA	-3	ASP	-	expression tag	UNP Q2KB35
AAA	-2	PRO	-	expression tag	UNP Q2KB35
AAA	-1	PHE	-	expression tag	UNP Q2KB35
AAA	0	THR	-	expression tag	UNP Q2KB35

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

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	BBB	1	Total	Zn	0	0
			1	1		
2	AAA	1	Total	Zn	0	0
			1	1		

- Molecule 3 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	BBB	2	Total Cl 2 2	0	0
3	AAA	3	Total Cl 3 3	0	0

- Molecule 4 is water.

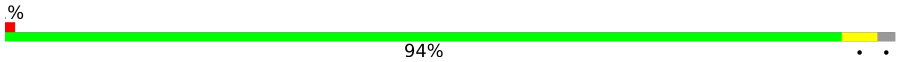
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	BBB	183	Total O 183 183	0	0
4	AAA	193	Total O 194 194	0	1

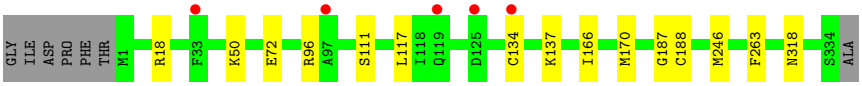
3 Residue-property plots (i)

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: Putative L-asparaginase II protein

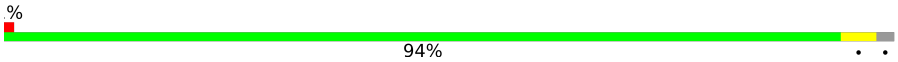
Chain BBB:





• Molecule 1: Putative L-asparaginase II protein

Chain AAA:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	96.25Å 85.30Å 82.95Å 90.00° 109.18° 90.00°	Depositor
Resolution (Å)	78.34 – 1.50 78.34 – 1.50	Depositor EDS
% Data completeness (in resolution range)	99.2 (78.34-1.50) 99.2 (78.34-1.50)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.09 (at 1.50Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.253 , 0.302 0.263 , 0.310	Depositor DCC
R_{free} test set	1000 reflections (1.00%)	wwPDB-VP
Wilson B-factor (Å^2)	19.4	Xtriage
Anisotropy	0.313	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.37 , 48.0	EDS
L-test for twinning[2]	$< \|L\| > = 0.45$, $< L^2 > = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	5337	wwPDB-VP
Average B, all atoms (Å^2)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.31% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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: ZN, CL, CSO

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	AAA	0.74	0/2510	0.79	0/3392
1	BBB	0.75	0/2518	0.79	0/3402
All	All	0.74	0/5028	0.79	0/6794

There are no bond length outliers.

There are no bond angle outliers.

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	AAA	2471	0	2451	10	0
1	BBB	2482	0	2462	14	0
2	AAA	1	0	0	0	0
2	BBB	1	0	0	0	0
3	AAA	3	0	0	1	0
3	BBB	2	0	0	0	0
4	AAA	194	0	0	4	0
4	BBB	183	0	0	4	0
All	All	5337	0	4913	22	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 2.

The worst 5 of 22 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:BBB:134[B]:CYS:HG	1:BBB:188[B]:CYS:HG	1.07	0.93
4:BBB:652:HOH:O	1:AAA:188[B]:CYS:SG	2.09	0.91
1:BBB:188[B]:CYS:SG	4:BBB:648:HOH:O	2.29	0.89
1:BBB:134[B]:CYS:CB	1:BBB:188[B]:CYS:HG	2.00	0.74
3:AAA:404:CL:CL	4:AAA:670[B]:HOH:O	2.50	0.66

There are no symmetry-related clashes.

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	AAA	332/341 (97%)	326 (98%)	6 (2%)	0	100	100
1	BBB	333/341 (98%)	325 (98%)	8 (2%)	0	100	100
All	All	665/682 (98%)	651 (98%)	14 (2%)	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	AAA	249/253 (98%)	248 (100%)	1 (0%)	91	82

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	BBB	250/253 (99%)	249 (100%)	1 (0%)	91 82
All	All	499/506 (99%)	497 (100%)	2 (0%)	91 82

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	BBB	263	PHE
1	AAA	263	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA (i)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains (i)

4 non-standard protein/DNA/RNA residues are modelled in this entry.

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	$\#\|Z\| > 2$
1	CSO	BBB	242[A]	-	3,6,7	0.78	0	0,6,8	-	-
1	CSO	AAA	242[A]	-	3,6,7	0.96	0	0,6,8	-	-
1	CSO	BBB	242[B]	-	3,6,7	1.10	0	0,6,8	-	-
1	CSO	AAA	242[B]	-	3,6,7	0.75	0	0,6,8	-	-

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
1	CSO	BBB	242[A]	-	-	0/1/5/7	-
1	CSO	AAA	242[A]	-	-	0/1/5/7	-
1	CSO	BBB	242[B]	-	-	0/1/5/7	-
1	CSO	AAA	242[B]	-	-	0/1/5/7	-

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.

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 7 ligands modelled in this entry, 7 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

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, 95^{th} 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	AAA	332/341 (97%)	0.29	3 (0%)	84	87	16, 23, 35, 46	0
1	BBB	333/341 (97%)	0.35	5 (1%)	73	78	16, 23, 34, 45	0
All	All	665/682 (97%)	0.32	8 (1%)	79	82	16, 23, 34, 46	0

The worst 5 of 8 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	BBB	33	PHE	4.3
1	AAA	2	THR	3.9
1	BBB	134[A]	CYS	2.7
1	AAA	334	SER	2.6
1	BBB	125	ASP	2.5

6.2 Non-standard residues in protein, DNA, RNA chains

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, 95^{th} 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
1	CSO	BBB	242[A]	7/8	0.94	0.14	16,18,19,21	4
1	CSO	BBB	242[B]	7/8	0.94	0.14	13,17,18,21	4
1	CSO	AAA	242[A]	7/8	0.95	0.12	13,16,17,19	4
1	CSO	AAA	242[B]	7/8	0.95	0.12	13,16,17,19	4

6.3 Carbohydrates

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, 95^{th} 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	CL	AAA	404	1/1	0.94	0.08	44,44,44,44	0
3	CL	BBB	403	1/1	0.96	0.04	41,41,41,41	0
3	CL	BBB	402	1/1	0.98	0.07	36,36,36,36	0
2	ZN	AAA	401	1/1	0.99	0.02	24,24,24,24	1
3	CL	AAA	402	1/1	0.99	0.07	40,40,40,40	0
3	CL	AAA	403	1/1	0.99	0.06	35,35,35,35	0
2	ZN	BBB	401	1/1	0.99	0.03	27,27,27,27	0

6.5 Other polymers (i)

There are no such residues in this entry.