

Full wwPDB X-ray Structure Validation Report (i)

Oct 6, 2023 – 05:32 AM EDT

PDB ID	:	8E9N
Title	:	Crystal structure of E. coli aspartate aminotransferase mutant VFIY in the
		ligand-free form at 278 K
Authors	:	Chica, R.A.; St-Jacques, A.D.; Rodriguez, J.M.; Thompson, M.C.
Deposited on		
Resolution	:	1.88 Å(reported)
Deposited on	:	Chica, R.A.; St-Jacques, A.D.; Rodriguez, J.M.; Thompson, M.C.

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 (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 (1)) were used in the production of this report:

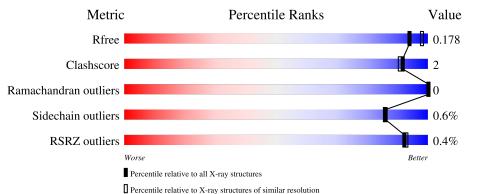
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.35.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35.1

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: $X\text{-}RAY \, DIFFRACTION$

The reported resolution of this entry is 1.88 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	$\begin{array}{c} \textbf{Whole archive} \\ (\#\textbf{Entries}) \end{array}$	${f Similar\ resolution}\ (\#{ m Entries,\ resolution\ range}({ m \AA}))$
R_{free}	130704	9470 (1.90-1.86)
Clashscore	141614	10282 (1.90-1.86)
Ramachandran outliers	138981	10152 (1.90-1.86)
Sidechain outliers	138945	10152 (1.90-1.86)
RSRZ outliers	127900	9303 (1.90-1.86)

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	А	406	93%	5% •					
1	В	406	95%						



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2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 12990 atoms, of which 6206 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 Aspartate aminotransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace	
1	А	399	Total 6356	C 2042	Н 3124	N 554	O 623	S 13	0	35	0
1	В	399	Total 6246	C 2014	Н 3066	N 544	O 609	S 13	0	27	0

Chain	Residue	Modelled	Actual		
А	-9	MET	-	initiating methionine	UNP P00509
А	-8	ALA	-	expression tag	UNP P00509
А	-7	HIS	-	expression tag	UNP P00509
А	-6	HIS	-	expression tag	UNP P00509
А	-5	HIS	-	expression tag	UNP P00509
A	-4	HIS	-	expression tag	UNP P00509
А	-3	HIS	-	expression tag	UNP P00509
А	-2	HIS	-	expression tag	UNP P00509
А	-1	VAL	-	expression tag	UNP P00509
А	0	GLY	-	expression tag	UNP P00509
А	1	THR	-	expression tag	UNP P00509
А	37	PHE	LYS	engineered mutation	UNP P00509
А	43	ILE	THR	engineered mutation	UNP P00509
А	64	TYR	ASN	engineered mutation	UNP P00509
В	-9	MET	-	initiating methionine	UNP P00509
В	-8	ALA	-	expression tag	UNP P00509
В	-7	HIS	-	expression tag	UNP P00509
В	-6	HIS	-	expression tag	UNP P00509
В	-5	HIS	-	expression tag	UNP P00509
В	-4	HIS	-	expression tag	UNP P00509
В	-3	HIS	-	expression tag	UNP P00509
В	-2	HIS	-	expression tag	UNP P00509
В	-1	VAL	-	expression tag	UNP P00509
В	0	GLY	-	expression tag	UNP P00509
В	1	THR	-	expression tag	UNP P00509

There are 28 discrepancies between the modelled and reference sequences:

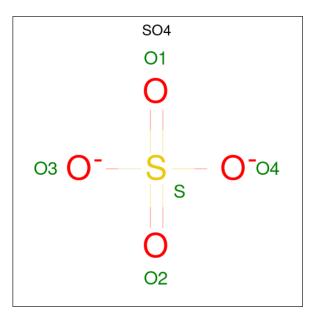
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Commu	Continued from previous page										
Chain	Residue	Modelled	Actual	Comment	Reference						
В	37	PHE	LYS	engineered mutation	UNP P00509						
В	43	ILE	THR	engineered mutation	UNP P00509						
В	64	TYR	ASN	engineered mutation	UNP P00509						

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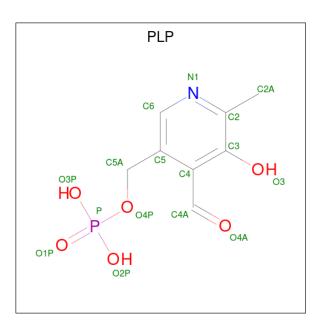
• Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	А	1	$\begin{array}{ccc} \text{Total} & \text{O} & \text{S} \\ 5 & 4 & 1 \end{array}$	0	0
2	В	1	$\begin{array}{ccc} \text{Total} & \text{O} & \text{S} \\ 5 & 4 & 1 \end{array}$	0	0

• Molecule 3 is PYRIDOXAL-5'-PHOSPHATE (three-letter code: PLP) (formula: $C_8H_{10}NO_6P$).





Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	
2	Δ	1	Total	С	Η	Ν	0	Р	0	0
0	D A	1	24	8	8	1	6	1	0	0
2	3 B	1	Total	С	Η	Ν	Ο	Р	0	0
J		1	24	8	8	1	6	1	0	0

• Molecule 4 is water.

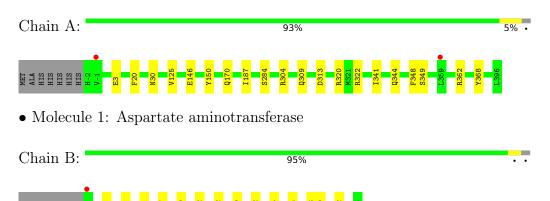
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	А	180	Total O 181 181	0	2
4	В	149	Total O 149 149	0	1



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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: Aspartate aminotransferase



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 63	Depositor
Cell constants	143.83Å 143.83Å 81.57Å	Depositor
a, b, c, α , β , γ	90.00° 90.00° 120.00°	Depositor
Resolution (Å)	124.43 - 1.88	Depositor
Itesolution (A)	124.56 - 1.88	EDS
% Data completeness	$98.6\ (124.43-1.88)$	Depositor
(in resolution range)	$98.6\ (124.56-1.88)$	EDS
R _{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	$1.07 (at 1.88 \text{\AA})$	Xtriage
Refinement program	PHENIX 1.17.1_3660	Depositor
R, R_{free}	0.150 , 0.178	Depositor
II, II, ree	0.150 , 0.178	DCC
R_{free} test set	7761 reflections (10.09%)	wwPDB-VP
Wilson B-factor $(Å^2)$	32.7	Xtriage
Anisotropy	0.161	Xtriage
Bulk solvent $k_{sol}(e/Å^3), B_{sol}(Å^2)$	0.38 , 40.5	EDS
L-test for twinning ²	$< L > = 0.50, < L^2 > = 0.33$	Xtriage
Estimated twinning fraction	0.037 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.98	EDS
Total number of atoms	12990	wwPDB-VP
Average B, all atoms $(Å^2)$	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 2.61% of the height of the origin peak. No significant pseudotranslation is detected.

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.



¹Intensities estimated from amplitudes.

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: PLP, $\mathrm{SO4}$

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		
	Unam	RMSZ	# Z > 5	RMSZ	# Z > 5	
1	А	0.33	0/3411	0.56	0/4626	
1	В	0.33	0/3327	0.54	0/4517	
All	All	0.33	0/6738	0.55	0/9143	

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 (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	А	3232	3124	3001	12	1
1	В	3180	3066	2984	9	1
2	А	5	0	0	0	0
2	В	5	0	0	0	0
3	А	16	8	8	0	0
3	В	16	8	8	0	0
4	А	181	0	0	3	0
4	В	149	0	0	1	0
All	All	6784	6206	6001	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 2.



Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:344:GLN:OE1	1:B:349[B]:SER:OG	1.89	0.89
1:A:344:GLN:OE1	1:A:349[B]:SER:OG	1.99	0.78
1:A:313[A]:ASP:OD1	4:A:501:HOH:O	2.08	0.71
1:A:309:GLN:NE2	4:A:501:HOH:O	2.19	0.66
1:B:322:ARG:NH1	1:B:348[A]:PHE:O	2.26	0.64
1:A:341[A]:ILE:HD12	1:A:349[A]:SER:HB2	1.87	0.56
1:A:322:ARG:NH1	1:A:348[A]:PHE:O	2.29	0.51
1:B:39[A]:GLU:HG2	1:B:382:PRO:CD	2.41	0.50
1:A:150:TYR:HE1	1:A:187:ILE:CD1	2.25	0.48
1:B:39[A]:GLU:HG2	1:B:382:PRO:HD2	1.96	0.47
1:A:362:ARG:HD2	1:A:368:TYR:CE2	2.51	0.45
1:B:344:GLN:OE1	1:B:349[B]:SER:CB	2.65	0.45
1:A:125[B]:VAL:O	1:A:146:GLU:HA	2.19	0.43
1:A:125[A]:VAL:O	1:A:146:GLU:HA	2.20	0.41
1:A:304:ARG:NH2	4:A:502:HOH:O	2.30	0.41
1:A:3:GLU:O	1:B:270:ARG:NH1	2.53	0.41
1:B:54:GLN:HG3	4:B:538:HOH:O	2.21	0.41
1:A:20:PHE:CZ	1:A:30:ASN:HB2	2.56	0.41
1:B:199:GLN:O	1:B:202:VAL:HG22	2.20	0.41
1:B:335:ASN:OD1	1:B:335:ASN:N	2.54	0.41

All (20) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

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)
1:A:320[B]:ARG:HH12	$1:B:24:GLU:OE2[2_545]$	1.52	0.08

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	Perce	ntiles	
1	А	432/406~(106%)	422 (98%)	10 (2%)	0	100	100
1	В	424/406 (104%)	414 (98%)	10 (2%)	0	100	100
All	All	856/812~(105%)	836~(98%)	20~(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 side chain 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	А	347/328~(106%)	344~(99%)	3(1%)	78 76
1	В	336/328~(102%)	333~(99%)	3 (1%)	78 76
All	All	683/656~(104%)	677~(99%)	6 (1%)	86 76

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

Mol	Chain	Res	Type
1	А	170[A]	GLN
1	А	170[B]	GLN
1	А	284	SER
1	В	167	ASN
1	В	212[A]	PHE
1	В	212[B]	PHE

Sometimes side chains can be flipped to improve hydrogen bonding and reduce clashes. There are no such side chains 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)

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)

4 ligands 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	Turne	Chain	Res	Link	Bond lengths			Bond angles		
IVIOI	Type	Chain	nes	LIIIK	Counts	RMSZ	# Z >2	Counts	RMSZ	# Z >2
3	PLP	В	402	-	16,16,16	4.54	9 (56%)	20,23,23	1.34	3 (15%)
2	SO4	А	401	-	4,4,4	0.14	0	6,6,6	0.33	0
3	PLP	А	402	-	16,16,16	4.43	9 (56%)	20,23,23	1.28	2 (10%)
2	SO4	В	401	-	4,4,4	0.14	0	6,6,6	0.31	0

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
3	PLP	В	402	-	-	2/8/8/8	0/1/1/1
3	PLP	А	402	-	-	3/8/8/8	0/1/1/1

All (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	В	402	PLP	C3-C2	11.48	1.52	1.40
3	А	402	PLP	C3-C2	11.04	1.51	1.40
3	В	402	PLP	C4-C3	-7.74	1.29	1.40
3	А	402	PLP	C4-C3	-7.68	1.29	1.40
3	В	402	PLP	C4-C5	6.78	1.50	1.42
3	В	402	PLP	C6-N1	6.64	1.48	1.34
3	А	402	PLP	C6-N1	6.56	1.48	1.34
3	А	402	PLP	C4-C5	6.44	1.50	1.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	В	402	PLP	C4-C4A	3.32	1.54	1.46
3	А	402	PLP	C2-N1	-3.22	1.27	1.33
3	А	402	PLP	C4-C4A	3.21	1.53	1.46
3	В	402	PLP	C2-N1	-3.03	1.28	1.33
3	А	402	PLP	C6-C5	-2.92	1.31	1.37
3	В	402	PLP	C6-C5	-2.91	1.31	1.37
3	В	402	PLP	P-O2P	-2.38	1.45	1.54
3	А	402	PLP	P-O3P	-2.30	1.46	1.54
3	А	402	PLP	P-O2P	-2.11	1.46	1.54
3	В	402	PLP	P-O3P	-2.09	1.46	1.54

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All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
3	В	402	PLP	O3P-P-O4P	3.24	115.35	106.73
3	А	402	PLP	O4P-P-O1P	2.94	114.72	106.47
3	В	402	PLP	C5-C6-N1	-2.37	119.87	123.82
3	В	402	PLP	O4P-P-O1P	2.32	112.98	106.47
3	А	402	PLP	C4-C3-C2	2.04	121.45	120.19

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	А	402	PLP	C5A-O4P-P-O2P
3	А	402	PLP	C5A-O4P-P-O3P
3	В	402	PLP	C5A-O4P-P-O2P
3	А	402	PLP	C3-C4-C4A-O4A
3	В	402	PLP	C3-C4-C4A-O4A

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, 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	$\langle RSRZ \rangle$	#RSRZ>2		$OWAB(Å^2)$	Q<0.9	
1	А	399/406~(98%)	-0.36	2 (0%)	91	91	22, 33, 57, 115	0
1	В	399/406~(98%)	-0.37	1 (0%)	94	94	22, 39, 65, 95	0
All	All	798/812~(98%)	-0.36	3 (0%)	92	93	22, 36, 60, 115	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	А	-1	VAL	5.4
1	А	359	LEU	2.4
1	В	-2	HIS	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 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	$\mathbf{B} ext{-factors}(\mathrm{\AA}^2)$	Q<0.9
2	SO4	А	401	5/5	0.99	0.09	$35,\!35,\!37,\!37$	0
2	SO4	В	401	5/5	0.99	0.12	34,36,36,37	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	$\mathbf{B} ext{-factors}(\mathrm{\AA}^2)$	Q < 0.9
3	PLP	А	402	16/16	0.99	0.11	$23,\!34,\!40,\!48$	0
3	PLP	В	402	16/16	0.99	0.10	24,42,50,58	0

6.5 Other polymers (i)

There are no such residues in this entry.

