



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 10, 2016 – 02:36 PM EDT

PDB ID : 1JAZ
Title : Crystal Structure of Monoclinic Form of D90E Mutant of Escherichia coli Asparaginase II
Authors : Borek, D.; Kozak, M.; Jaskolski, M.
Deposited on : 2001-06-01
Resolution : 2.27 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

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	unknown
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20027939
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	rb-20027939

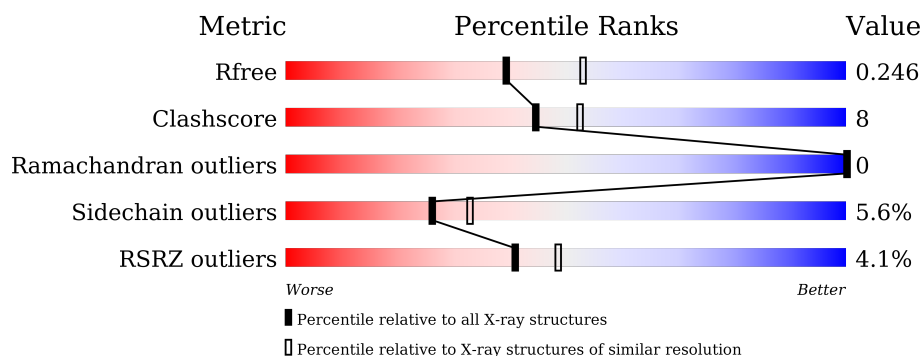
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.27 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}	91344	5193 (2.30-2.26)
Clashscore	102246	5929 (2.30-2.26)
Ramachandran outliers	100387	5851 (2.30-2.26)
Sidechain outliers	100360	5850 (2.30-2.26)
RSRZ outliers	91569	5204 (2.30-2.26)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	326	<div> <div>3%</div> <div> <div></div> <div>73%</div> <div>19%</div> <div>• 6%</div> </div> </div>
1	B	326	<div> <div>4%</div> <div> <div></div> <div>76%</div> <div>16%</div> <div>• 6%</div> </div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4739 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 L-ASPARAGINASE II.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	307	Total	C	N	O	S	0	2	0
			2309	1444	395	462	8			
1	B	305	Total	C	N	O	S	0	1	0
			2291	1432	392	459	8			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	90	GLU	ASP	ENGINEERED	UNP P00805
B	90	GLU	ASP	ENGINEERED	UNP P00805

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

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	2	Total	Zn	0	0
			2	2		
2	A	1	Total	Zn	0	0
			1	1		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	71	Total	O	0	0
			71	71		
3	B	65	Total	O	0	0
			65	65		

- Molecule 1: L-ASPARAGINASE II



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	73.12Å 133.08Å 62.56Å 90.00° 108.78° 90.00°	Depositor
Resolution (Å)	10.00 – 2.27 10.00 – 2.27	Depositor EDS
% Data completeness (in resolution range)	99.5 (10.00-2.27) 99.5 (10.00-2.27)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.43 (at 2.26Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.180 , 0.232 0.197 , 0.246	Depositor DCC
R_{free} test set	1301 reflections (5.33%)	DCC
Wilson B-factor (Å ²)	33.3	Xtriage
Anisotropy	0.049	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.44 , 47.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4739	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 6.65% 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

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	0.74	0/2355	0.97	8/3207 (0.2%)
1	B	0.76	0/2332	0.97	11/3175 (0.3%)
All	All	0.75	0/4687	0.97	19/6382 (0.3%)

There are no bond length outliers.

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	315	ASP	CB-CG-OD2	6.94	124.55	118.30
1	A	44	ASP	CB-CG-OD2	6.75	124.38	118.30
1	B	286	ASP	CB-CG-OD2	6.74	124.37	118.30
1	B	156	ASP	CB-CG-OD2	6.20	123.88	118.30
1	B	60	ASP	CB-CG-OD2	6.14	123.83	118.30
1	B	106	ASP	CB-CG-OD2	6.02	123.72	118.30
1	B	159	ASP	CB-CG-OD2	5.97	123.67	118.30
1	A	76	ASP	CB-CG-OD2	5.87	123.58	118.30
1	A	286	ASP	CB-CG-OD2	5.75	123.48	118.30
1	A	225	ASP	CB-CG-OD2	5.72	123.45	118.30
1	B	233	ASP	CB-CG-OD2	5.68	123.41	118.30
1	B	76	ASP	CB-CG-OD2	5.49	123.24	118.30
1	B	167	ASP	CB-CG-OD1	5.38	123.14	118.30
1	A	159	ASP	CB-CG-OD2	5.27	123.05	118.30
1	B	81	ASP	CB-CG-OD2	5.26	123.03	118.30
1	A	285	ASP	CB-CG-OD2	5.21	122.99	118.30
1	B	315	ASP	CB-CG-OD2	5.15	122.93	118.30
1	A	156	ASP	CB-CG-OD2	5.07	122.86	118.30
1	B	100	ASP	CB-CG-OD2	5.04	122.84	118.30

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	2309	0	2305	43	0
1	B	2291	0	2283	31	0
2	A	1	0	0	0	0
2	B	2	0	0	0	0
3	A	71	0	0	0	0
3	B	65	0	0	5	0
All	All	4739	0	4588	73	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 (73) 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:248:ASN:HD22	1:A:249:LEU:H	1.25	0.85
1:A:300:GLN:HE21	1:A:300:GLN:H	1.25	0.81
1:B:39:VAL:HG22	1:B:42:LEU:HG	1.67	0.76
1:A:246:ASN:HB3	1:A:283:GLU:HG3	1.67	0.76
1:A:58:SER:OG	1:A:90:GLU:HG2	1.86	0.75
1:A:300:GLN:NE2	1:A:300:GLN:H	1.86	0.73
1:A:64:ASN:O	1:A:68:THR:HG23	1.89	0.72
1:B:164:ASN:HD22	1:B:167:ASP:H	1.41	0.68
1:A:196:LYS:NZ	1:A:325:GLN:NE2	2.43	0.67
1:A:300:GLN:N	1:A:300:GLN:HE21	1.92	0.66
1:A:40:PRO:O	1:A:43:LYS:HG2	1.97	0.64
1:B:281:ASP:O	1:B:281:ASP:CG	2.34	0.64
1:A:196:LYS:NZ	1:A:325:GLN:HE21	1.97	0.63
1:A:39:VAL:O	1:A:42:LEU:HB2	2.00	0.62
1:B:64:ASN:O	1:B:68:THR:HG23	2.01	0.60
1:A:158:ARG:CZ	1:A:297:LEU:HD21	2.32	0.60
1:B:281:ASP:O	1:B:281:ASP:OD1	2.21	0.58
1:A:246:ASN:HB3	1:A:283:GLU:CG	2.34	0.57
1:B:13:ILE:O	1:B:125:GLY:HA3	2.05	0.57
1:B:201:THR:O	3:B:427:HOH:O	2.18	0.56
1:A:310:LEU:HD23	1:A:314:LYS:HD3	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:163:THR:CG2	3:B:435:HOH:O	2.55	0.53
1:B:251:LYS:HG2	1:B:252:SER:N	2.23	0.53
1:B:163:THR:HG21	3:B:435:HOH:O	2.08	0.53
1:A:183[A]:HIS:NE2	1:A:184:ASN:ND2	2.56	0.53
1:A:36:VAL:O	1:A:40:PRO:HA	2.09	0.53
1:A:298:ASN:HB2	1:A:299:PRO:CD	2.39	0.53
1:A:248:ASN:ND2	1:A:249:LEU:H	2.02	0.52
1:B:281:ASP:C	1:B:281:ASP:OD1	2.48	0.52
1:B:39:VAL:O	1:B:39:VAL:HG22	2.09	0.52
1:A:248:ASN:HD22	1:A:249:LEU:N	2.02	0.51
1:B:163:THR:HB	1:B:170:THR:O	2.11	0.51
1:B:58:SER:OG	1:B:90:GLU:HB2	2.11	0.50
1:A:110:VAL:HG13	1:A:149:VAL:HG23	1.93	0.50
1:A:196:LYS:HZ1	1:A:325:GLN:NE2	2.09	0.50
1:A:90:GLU:CD	1:A:90:GLU:H	2.12	0.50
1:A:196:LYS:HE3	1:A:325:GLN:HB3	1.94	0.49
1:B:182:ILE:HG12	1:B:187:ILE:HG12	1.95	0.49
1:A:8:ALA:HA	1:A:86:THR:OG1	2.13	0.49
1:A:196:LYS:HZ2	1:A:325:GLN:NE2	2.11	0.48
1:B:248:ASN:HD22	1:B:249:LEU:H	1.59	0.48
1:A:39:VAL:O	1:A:39:VAL:HG22	2.13	0.48
1:A:157:GLY:O	1:A:301:LYS:NZ	2.47	0.48
1:B:73:ILE:HD11	1:B:85:ILE:HD11	1.95	0.48
1:B:42:LEU:HD11	1:B:129:LEU:CD1	2.45	0.47
1:A:160:VAL:O	1:A:301:LYS:HE3	2.15	0.47
1:B:183[A]:HIS:CD2	1:B:184:ASN:HB2	2.50	0.46
1:A:6:ILE:HD12	1:A:48:VAL:HG11	1.98	0.45
1:A:196:LYS:CE	1:A:325:GLN:HE21	2.29	0.44
1:B:246:ASN:HB3	1:B:283:GLU:HG3	1.99	0.44
1:A:56:ILE:O	1:A:56:ILE:HG13	2.17	0.44
1:A:13:ILE:O	1:A:125:GLY:HA3	2.18	0.43
1:A:12:THR:OG1	1:A:14:ALA:O	2.26	0.43
1:A:56:ILE:HD11	1:A:61:MET:HE2	2.01	0.43
1:B:8:ALA:HA	1:B:86:THR:OG1	2.18	0.43
1:B:56:ILE:HD11	1:B:61:MET:HE2	2.01	0.43
1:B:270:SER:HA	1:B:294:SER:OG	2.18	0.43
1:A:53:VAL:HG12	1:A:54:VAL:HG13	2.01	0.43
1:A:315:ASP:OD2	1:A:317:GLN:HB2	2.18	0.42
1:B:229:LYS:NZ	3:B:438:HOH:O	2.52	0.42
1:A:66:TRP:HB3	1:A:98:PHE:CE2	2.54	0.42
1:B:298:ASN:HB2	1:B:299:PRO:CD	2.49	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:214:VAL:HA	1:B:238:GLY:O	2.19	0.42
1:B:53:VAL:HG12	1:B:54:VAL:HG13	2.01	0.41
1:A:216:ILE:HG12	1:A:240:VAL:HB	2.01	0.41
1:A:196:LYS:HE3	1:A:325:GLN:HE21	1.86	0.41
1:B:248:ASN:C	1:B:249:LEU:HG	2.36	0.41
1:A:3:ASN:OD1	1:A:3:ASN:O	2.38	0.41
1:B:298:ASN:HA	3:B:410:HOH:O	2.21	0.41
1:A:122:SER:HB2	1:B:127:PHE:HB2	2.03	0.41
1:B:39:VAL:N	1:B:40:PRO:HD3	2.35	0.41
1:A:310:LEU:O	1:A:314:LYS:HE3	2.22	0.40

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	A	305/326 (94%)	297 (97%)	8 (3%)	0	100	100
1	B	302/326 (93%)	295 (98%)	7 (2%)	0	100	100
All	All	607/652 (93%)	592 (98%)	15 (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	A	254/266 (96%)	237 (93%)	17 (7%)	20	24
1	B	251/266 (94%)	240 (96%)	11 (4%)	35	45
All	All	505/532 (95%)	477 (94%)	28 (6%)	26	34

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	37	ASN
1	A	39	VAL
1	A	41	GLN
1	A	68	THR
1	A	129	LEU
1	A	143	ASN
1	A	156	ASP
1	A	179	LEU
1	A	197	HIS
1	A	211	LEU
1	A	222	ASN
1	A	248	ASN
1	A	251	LYS
1	A	254	PHE
1	A	257	LEU
1	A	300	GLN
1	A	308	LEU
1	B	12	THR
1	B	39	VAL
1	B	41	GLN
1	B	156	ASP
1	B	163	THR
1	B	197	HIS
1	B	222	ASN
1	B	248	ASN
1	B	254	PHE
1	B	281	ASP
1	B	308	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	131	ASN
1	A	184	ASN

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Mol	Chain	Res	Type
1	A	248	ASN
1	A	280	GLN
1	A	300	GLN
1	A	318	GLN
1	A	325	GLN
1	B	164	ASN
1	B	248	ASN
1	B	318	GLN
1	B	325	GLN

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 3 ligands modelled in this entry, 3 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 ⓘ

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, 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	307/326 (94%)	-0.02	11 (3%) 46 54	17, 23, 32, 44	0
1	B	305/326 (93%)	-0.02	14 (4%) 36 44	17, 22, 34, 42	0
All	All	612/652 (93%)	-0.02	25 (4%) 41 48	17, 23, 34, 44	0

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	281	ASP	5.4
1	A	37	ASN	3.6
1	A	281	ASP	3.4
1	B	11	GLY	3.3
1	B	44	ASP	2.9
1	B	287	ALA	2.8
1	B	78	ASP	2.8
1	B	3	ASN	2.7
1	B	40	PRO	2.7
1	B	317	GLN	2.6
1	B	184	ASN	2.6
1	A	317	GLN	2.6
1	A	78	ASP	2.5
1	B	262	LYS	2.5
1	A	321	GLN	2.5
1	B	38	ALA	2.5
1	B	321	GLN	2.4
1	A	3	ASN	2.3
1	B	106	ASP	2.3
1	A	36	VAL	2.2
1	B	41	GLN	2.1
1	A	209	ASN	2.1
1	A	11	GLY	2.1
1	A	15	GLY	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	233	ASP	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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
2	ZN	A	401	1/1	0.98	0.06	-2.19	37,37,37,37	0
2	ZN	B	402	1/1	0.99	0.07	-7.08	33,33,33,33	0
2	ZN	B	403	1/1	0.61	0.13	-	43,43,43,43	1

6.5 Other polymers [i](#)

There are no such residues in this entry.