



# wwPDB X-ray Structure Validation Summary Report i

Feb 28, 2014 – 06:52 PM GMT

PDB ID : 2AQO  
Title : Crystal structure of E. coli Isoaspartyl Dipeptidase Mutant E77Q  
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Deposited on : 2005-08-18  
Resolution : 1.95 Å(reported)

This is a wwPDB validation summary report for a publicly released PDB entry.

We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)

A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

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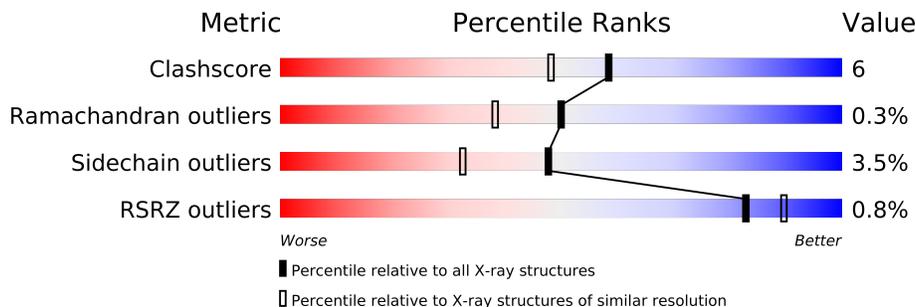
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : dev-1323  
EDS : stable22639  
Percentile statistics : 21963  
Refmac : 5.8.0049  
CCP4 : 6.3.0 (Settle)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance

The reported resolution of this entry is 1.95 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))
Clashscore	79885	1488 (1.96-1.96)
Ramachandran outliers	78287	1475 (1.96-1.96)
Sidechain outliers	78261	1475 (1.96-1.96)
RSRZ outliers	66119	1321 (1.96-1.96)

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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	390	
1	B	390	

## 2 Entry composition i

There are 3 unique types of molecules in this entry. The entry contains 6005 atoms, of which 0 are hydrogen and 0 are deuterium.

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 Isoaspartyl dipeptidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	378	2799	1762	482	543	12	0	2	0
1	B	375	2780	1750	479	539	12	0	2	0

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	77	GLN	GLU	ENGINEERED	UNP P39377
A	162	KCX	LYS	MODIFIED RESIDUE	UNP P39377
B	77	GLN	GLU	ENGINEERED	UNP P39377
B	162	KCX	LYS	MODIFIED RESIDUE	UNP P39377

- 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	2	Total	Zn	0	0
			2	2		

- Molecule 3 is water.

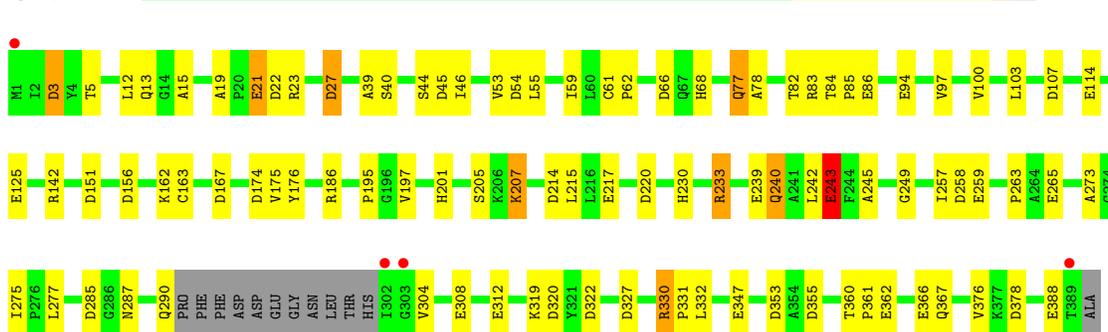
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	207	Total	O	0	0
			207	207		
3	B	215	Total	O	0	0
			215	215		

### 3 Residue-property plots i

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

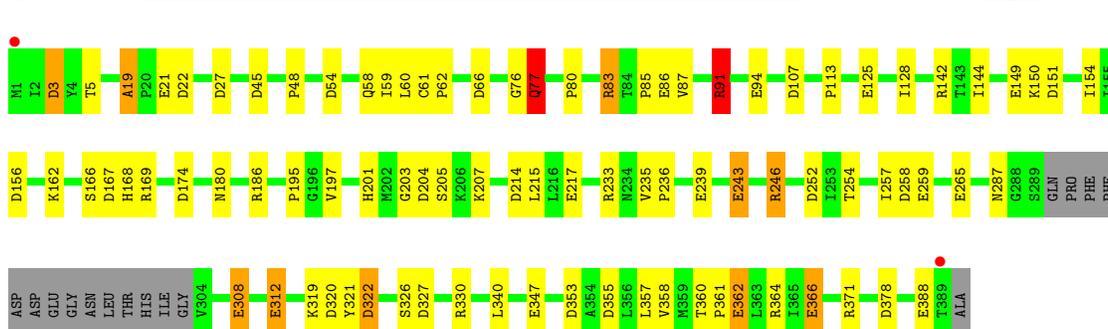
- Molecule 1: Isoaspartyl dipeptidase

Chain A:



- Molecule 1: Isoaspartyl dipeptidase

Chain B:



## 4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 4 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	119.20Å 119.20Å 138.60Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 1.95 36.51 – 1.93	Depositor EDS
% Data completeness (in resolution range)	93.4 (30.00-1.95) 91.2 (36.51-1.93)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.07	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.49 (at 1.94Å)	Xtrriage
Refinement program	TNT V. 5-E	Depositor
R, $R_{free}$	0.179 , 0.219 0.159 , (Not available)	Depositor DCC
$R_{free}$ test set	No test flags present.	DCC
Wilson B-factor (Å <sup>2</sup> )	20.0	Xtrriage
Anisotropy	0.196	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 68.6	EDS
Estimated twinning fraction	No twinning to report.	Xtrriage
L-test for twinning	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtrriage
Outliers	1 of 68348 reflections (0.001%)	Xtrriage
$F_o, F_c$ correlation	0.97	EDS
Total number of atoms	6005	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	29.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.96% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities 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: ZN, KCX

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	0.94	17/2839 (0.6%)	1.38	37/3861 (1.0%)
1	B	0.95	16/2820 (0.6%)	1.39	44/3834 (1.1%)
All	All	0.95	33/5659 (0.6%)	1.39	81/7695 (1.1%)

The worst 5 of 33 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	243	GLU	CD-OE2	8.65	1.35	1.25
1	A	114	GLU	CD-OE2	8.03	1.34	1.25
1	B	149	GLU	CD-OE2	7.66	1.34	1.25
1	B	259	GLU	CD-OE2	7.03	1.33	1.25
1	A	217	GLU	CD-OE2	7.02	1.33	1.25

The worst 5 of 81 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
1	A	233	ARG	NE-CZ-NH1	10.64	125.62	120.30
1	B	66	ASP	CB-CG-OD2	-10.40	108.94	118.30
1	A	233	ARG	NE-CZ-NH2	-9.23	115.68	120.30
1	A	83	ARG	NE-CZ-NH1	9.23	124.91	120.30
1	B	22	ASP	CB-CG-OD1	8.72	126.15	118.30

There are no chirality outliers.

There are no planarity outliers.

### 5.2 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 hydrogens

added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2799	0	2839	38	0
1	B	2780	0	2821	33	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
3	A	207	0	0	3	0
3	B	215	0	0	5	2
All	All	6005	0	5660	70	2

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 6.

The worst 5 of 70 close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:59:ILE:HD12	1:B:361:PRO:HA	1.70	0.72
1:B:205:SER:OG	1:B:207:LYS:HG2	1.94	0.66
1:A:142:ARG:HB2	3:A:950:HOH:O	1.96	0.66
1:A:59:ILE:CD1	1:A:361:PRO:HA	2.27	0.65
1:B:83:ARG:HG3	3:B:811:HOH:O	1.96	0.65

All (2) 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	Distance(Å)	Clash(Å)
3:B:915:HOH:O	3:B:915:HOH:O[7_556]	1.75	0.45
3:B:913:HOH:O	3:B:913:HOH:O[7_556]	1.98	0.22

## 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	375/390 (96%)	365 (97%)	9 (2%)	1 (0%)	50 38

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	372/390 (95%)	365 (98%)	6 (2%)	1 (0%)	50	38
All	All	747/780 (96%)	730 (98%)	15 (2%)	2 (0%)	50	38

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	19	ALA
1	B	19	ALA

### 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	304/312 (97%)	292 (96%)	12 (4%)	43	27
1	B	302/312 (97%)	293 (97%)	9 (3%)	53	39
All	All	606/624 (97%)	585 (96%)	21 (4%)	48	32

5 of 21 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	290	GLN
1	A	319	LYS
1	B	319	LYS
1	A	243	GLU
1	B	326	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 10 such sidechains are listed below:

Mol	Chain	Res	Type
1	A	367	GLN
1	B	13	GLN
1	B	168	HIS
1	A	290	GLN
1	B	77	GLN

### 5.3.3 RNA ⓘ

There are no RNA chains in this entry.

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

2 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	KCX	A	162	1,2	11,11,12	6.73	3 (27%)	10,12,14	4.17	4 (40%)
1	KCX	B	162	1,2	11,11,12	7.62	3 (27%)	10,12,14	2.94	4 (40%)

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	KCX	A	162	1,2	-	0/8/10/12	0/0/0/0
1	KCX	B	162	1,2	-	0/8/10/12	0/0/0/0

The worst 5 of 6 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	162	KCX	O-C	19.95	1.25	1.11
1	A	162	KCX	O-C	17.27	1.23	1.11
1	B	162	KCX	CX-NZ	15.26	1.47	1.32
1	A	162	KCX	CX-NZ	13.67	1.45	1.32
1	A	162	KCX	CA-C	3.14	1.54	1.48

The worst 5 of 8 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	162	KCX	CE-NZ-CX	-10.93	102.08	121.99
1	B	162	KCX	CE-NZ-CX	-5.78	111.46	121.99

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	162	KCX	C-CA-N	-5.21	108.63	113.83
1	B	162	KCX	C-CA-N	-5.19	108.64	113.83
1	A	162	KCX	OQ2-CX-NZ	-4.07	111.26	116.33

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

## 5.5 Carbohydrates

There are no carbohydrates in this entry.

## 5.6 Ligand geometry

Of 4 ligands modelled in this entry, 4 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.

## 5.7 Other polymers

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	378/390 (96%)	-0.49	4 (1%) 77 85	13, 24, 59, 100	0
1	B	375/390 (96%)	-0.54	2 (0%) 88 94	13, 23, 55, 99	0
All	All	753/780 (96%)	-0.52	6 (0%) 83 89	13, 24, 56, 100	0

The worst 5 of 6 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	389	THR	3.7
1	A	389	THR	3.6
1	B	1	MET	3.4
1	A	303	GLY	2.8
1	A	302	ILE	2.7

### 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. 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, 95<sup>th</sup> 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	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
1	KCX	B	162	12/13	0.10	0.96	11,16,31,35	0
1	KCX	A	162	12/13	0.10	0.80	14,20,29,43	0

### 6.3 Carbohydrates

There are no carbohydrates in this entry.

## 6.4 Ligands

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, 95<sup>th</sup> 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	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
2	ZN	A	801	1/1	0.03	-1.85	35,35,35,35	0
2	ZN	A	802	1/1	0.03	-2.37	38,38,38,38	0
2	ZN	B	803	1/1	0.02	-2.76	32,32,32,32	0
2	ZN	B	804	1/1	0.03	-3.65	36,36,36,36	0

## 6.5 Other polymers

There are no such residues in this entry.