



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 05:18 AM GMT

PDB ID : 2Q7S  
Title : Crystal structure of N-formylglutamate amidohydrolase (YP\_297560.1) from *Ralstonia eutropha* JMP134 at 2.00 Å resolution  
Authors : Joint Center for Structural Genomics (JCSG)  
Deposited on : 2007-06-07  
Resolution : 2.00 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](mailto: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.

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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 : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
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) : trunk26865

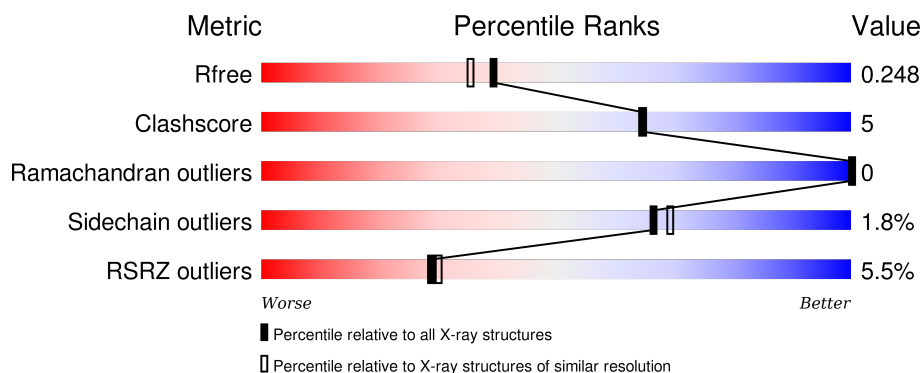
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.00 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	6249 (2.00-2.00)
Clashscore	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (2.00-2.00)
RSRZ outliers	91569	6262 (2.00-2.00)

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. 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	290	<div> <div>6%</div> <div> <div></div> <div>81%</div> <div>12%</div> <div>7%</div> </div> </div>
1	B	290	<div> <div>4%</div> <div> <div></div> <div>85%</div> <div>11%</div> <div>•</div> </div> </div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4703 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 N-formylglutamate amidohydrolase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	271	Total	C	N	O	S	Se	0	2	0
			2153	1366	388	394	2	3			
1	B	280	Total	C	N	O	S	Se	0	8	0
			2251	1434	402	409	2	4			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	GLY	-	LEADER SEQUENCE	UNP Q46VW8
A	1	MSE	MET	MODIFIED RESIDUE	UNP Q46VW8
A	5	MSE	MET	MODIFIED RESIDUE	UNP Q46VW8
A	86	MSE	MET	MODIFIED RESIDUE	UNP Q46VW8
A	173	MSE	MET	MODIFIED RESIDUE	UNP Q46VW8
A	256	MSE	MET	MODIFIED RESIDUE	UNP Q46VW8
B	0	GLY	-	LEADER SEQUENCE	UNP Q46VW8
B	1	MSE	MET	MODIFIED RESIDUE	UNP Q46VW8
B	5	MSE	MET	MODIFIED RESIDUE	UNP Q46VW8
B	86	MSE	MET	MODIFIED RESIDUE	UNP Q46VW8
B	173	MSE	MET	MODIFIED RESIDUE	UNP Q46VW8
B	256	MSE	MET	MODIFIED RESIDUE	UNP Q46VW8

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

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

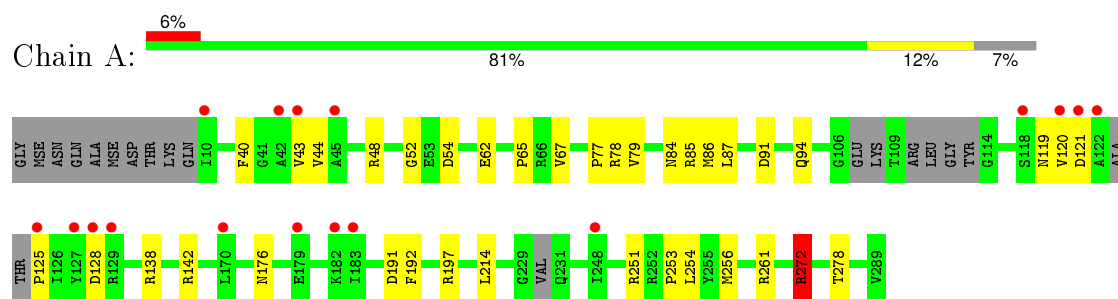
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	152	Total 152	O 152	0	0
3	B	145	Total 145	O 145	0	0

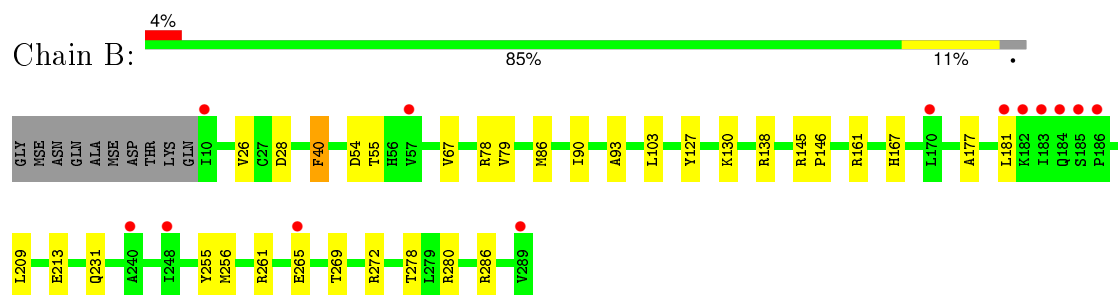
### 3 Residue-property plots [i](#)

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: N-formylglutamate amidohydrolase



- Molecule 1: N-formylglutamate amidohydrolase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	58.61Å 74.83Å 145.16Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.60 – 2.00 29.60 – 2.00	Depositor EDS
% Data completeness (in resolution range)	99.8 (29.60-2.00) 99.8 (29.60-2.00)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	0.08	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.49 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, $R_{free}$	0.199 , 0.246 0.203 , 0.248	Depositor DCC
$R_{free}$ test set	2213 reflections (5.31%)	DCC
Wilson B-factor (Å <sup>2</sup> )	26.9	Xtriage
Anisotropy	0.122	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 48.5	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	5 of 43899 reflections (0.011%)	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	4703	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 36.89 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 4.6607e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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	0.82	0/2206	0.85	4/2997 (0.1%)
1	B	0.80	0/2327	0.82	5/3166 (0.2%)
All	All	0.81	0/4533	0.83	9/6163 (0.1%)

There are no bond length outliers.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	272	ARG	NE-CZ-NH1	7.00	123.80	120.30
1	A	54	ASP	CB-CG-OD2	-5.47	113.38	118.30
1	B	161	ARG	NE-CZ-NH2	-5.28	117.66	120.30
1	A	85	ARG	NE-CZ-NH1	5.21	122.90	120.30
1	B	138	ARG	NE-CZ-NH1	5.14	122.87	120.30
1	B	54	ASP	CB-CG-OD2	-5.08	113.73	118.30
1	B	54	ASP	CB-CG-OD1	5.05	122.84	118.30
1	B	286	ARG	NE-CZ-NH1	5.05	122.82	120.30
1	A	197	ARG	NE-CZ-NH2	-5.03	117.78	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	2153	0	2126	21	0
1	B	2251	0	2245	24	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	152	0	0	5	0
3	B	145	0	0	3	0
All	All	4703	0	4371	45	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 5.

All (45) 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:B:177:ALA:O	1:B:181:LEU:HD13	1.72	0.88
1:A:86:MSE:HE2	3:A:458:HOH:O	1.80	0.80
1:A:256:MSE:HE1	1:A:261:ARG:CZ	2.15	0.75
1:B:55:THR:HB	1:B:256[B]:MSE:HE1	1.68	0.74
1:A:119:ASN:HB3	1:A:125:PRO:HA	1.72	0.72
1:B:256[B]:MSE:HE2	1:B:261:ARG:HA	1.73	0.71
1:A:43:VAL:HG11	1:A:128:ASP:HB3	1.76	0.68
1:B:209[A]:LEU:HD11	1:B:280[A]:ARG:HG3	1.76	0.67
1:B:209[A]:LEU:HD11	1:B:280[A]:ARG:CG	2.28	0.63
1:B:67:VAL:HG21	1:B:278:THR:HG22	1.82	0.61
1:A:86:MSE:CE	3:A:458:HOH:O	2.45	0.60
1:A:272:ARG:HG3	1:A:272:ARG:HH11	1.64	0.60
1:B:269:THR:HG23	1:B:272:ARG:HH12	1.65	0.60
1:B:269:THR:HG23	1:B:272:ARG:NH1	2.17	0.60
1:B:256[B]:MSE:HE2	1:B:261:ARG:CA	2.34	0.57
1:A:191:ASP:HB3	1:A:254:LEU:HD11	1.88	0.56
1:A:67:VAL:HG21	1:A:278:THR:HG22	1.88	0.53
1:A:44:VAL:HG22	1:A:48:ARG:HB2	1.90	0.53
1:B:265:GLU:CD	1:B:265:GLU:H	2.12	0.53
1:B:255:TYR:CE2	1:B:256[A]:MSE:HG2	2.44	0.52
1:A:251:ARG:HG2	1:A:253:PRO:HD2	1.92	0.51
1:B:55:THR:CB	1:B:256[B]:MSE:HE1	2.38	0.49
1:B:103:LEU:HB2	3:B:519:HOH:O	2.12	0.48
1:B:40:PHE:HE1	1:B:127:TYR:HH	1.59	0.48
1:A:52:GLY:HA2	3:A:523:HOH:O	2.13	0.48
1:B:280[B]:ARG:HG2	3:B:472:HOH:O	2.13	0.47
1:A:272:ARG:HH11	1:A:272:ARG:CG	2.26	0.47
1:B:93:ALA:O	1:B:130:LYS:HE2	2.14	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:91:ASP:O	1:A:94:GLN:HG2	2.16	0.46
1:A:192:PHE:CG	1:A:214:LEU:HD13	2.51	0.46
1:A:120:VAL:HG23	1:A:121:ASP:OD1	2.16	0.45
1:B:145:ARG:HB2	1:B:146:PRO:HD3	1.97	0.45
1:A:138:ARG:HD2	1:A:142:ARG:NH2	2.33	0.44
1:B:209[B]:LEU:HD22	1:B:213:GLU:HG2	2.00	0.43
1:B:26:VAL:O	1:B:167:HIS:HA	2.19	0.43
1:A:176:ASN:HB3	3:A:513:HOH:O	2.18	0.43
1:A:62:GLU:O	1:A:65:PRO:HD2	2.18	0.42
1:B:90:ILE:HB	1:B:103:LEU:HD22	2.01	0.41
1:B:86:MSE:HG3	1:B:231[B]:GLN:OE1	2.18	0.41
1:A:87:LEU:HA	1:A:87:LEU:HD23	1.83	0.41
1:A:86:MSE:HE3	3:A:406:HOH:O	2.20	0.41
1:B:256[A]:MSE:HE1	1:B:261:ARG:CZ	2.50	0.40
1:B:255:TYR:CD2	1:B:256[A]:MSE:HG2	2.57	0.40
1:B:79[B]:VAL:HG13	3:B:417:HOH:O	2.21	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	264/290 (91%)	260 (98%)	4 (2%)	0	100	100
1	B	286/290 (99%)	281 (98%)	5 (2%)	0	100	100
All	All	550/580 (95%)	541 (98%)	9 (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	223/234 (95%)	218 (98%)	5 (2%)	60	62
1	B	233/234 (100%)	230 (99%)	3 (1%)	76	79
All	All	456/468 (97%)	448 (98%)	8 (2%)	66	69

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	40	PHE
1	A	77	PRO
1	A	78	ARG
1	A	84	ASN
1	A	272	ARG
1	B	28	ASP
1	B	40	PHE
1	B	78	ARG

Some 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](#)

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

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

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	268/290 (92%)	0.30	17 (6%) 23 24	10, 17, 34, 44	0
1	B	277/290 (95%)	0.18	13 (4%) 35 37	10, 18, 31, 44	0
All	All	545/580 (93%)	0.24	30 (5%) 29 30	10, 18, 33, 44	0

All (30) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	184	GLN	4.9
1	A	10	ILE	4.4
1	A	118	SER	3.9
1	A	125	PRO	3.5
1	A	45	ALA	3.4
1	A	248	ILE	3.4
1	B	10	ILE	3.3
1	B	186	PRO	3.1
1	A	182	LYS	3.0
1	B	57	VAL	2.8
1	A	128	ASP	2.8
1	A	42	ALA	2.8
1	A	170	LEU	2.8
1	A	120	VAL	2.8
1	B	240	ALA	2.7
1	A	129	ARG	2.7
1	A	121	ASP	2.6
1	B	170	LEU	2.6
1	B	248	ILE	2.6
1	B	185	SER	2.6
1	A	122	ALA	2.5
1	A	127	TYR	2.5
1	B	181	LEU	2.5
1	A	43	VAL	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	179	GLU	2.4
1	A	183	ILE	2.3
1	B	265	GLU	2.3
1	B	183	ILE	2.3
1	B	182	LYS	2.1
1	B	289	VAL	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, 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	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
2	ZN	A	400	1/1	0.97	0.09	-3.68	27,27,27,27	0
2	ZN	B	400	1/1	0.99	0.08	-9.80	25,25,25,25	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.