

# Full wwPDB X-ray Structure Validation Report (i)

#### Nov 1, 2023 - 01:09 PM JST

PDB ID : 5HXD

Title : Crystal structure of murein-tripeptide amidase MpaA from Escherichia coli

O157

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Χ.

Deposited on : 2016-01-30

Resolution : 2.60 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
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:

Mol Probity : 4.02b-467

Mogul: 1.8.5 (274361), 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)

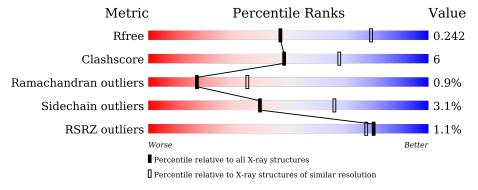
Validation Pipeline (wwPDB-VP) : 2.36

### 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 2.60 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	$\begin{array}{c} \textbf{Whole archive} \\ (\#\text{Entries}) \end{array}$	$\begin{array}{c} {\rm Similar\ resolution} \\ (\#{\rm Entries},\ {\rm resolution\ range}({\rm \AA})) \end{array}$
$R_{free}$	130704	3163 (2.60-2.60)
Clashscore	141614	3518 (2.60-2.60)
Ramachandran outliers	138981	3455 (2.60-2.60)
Sidechain outliers	138945	3455 (2.60-2.60)
RSRZ outliers	127900	3104 (2.60-2.60)

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	A	237	82%	17%
1	В	237	83%	16% •



### 2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 3700 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.

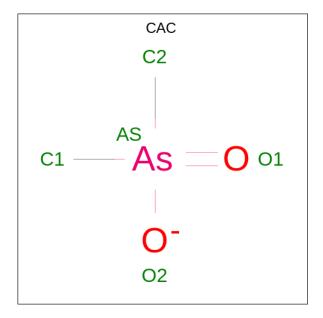
• Molecule 1 is a protein called Protein MpaA.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	Λ	237	Total	С	Ν	О	S	0	0	0
1	A	231	1834	1157	326	342	9	U	U	
1	D	234	Total	С	N	О	S	0	0	0
1	Ъ	204	1817	1147	323	338	9		U	

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

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

• Molecule 3 is CACODYLATE ION (three-letter code: CAC) (formula: C<sub>2</sub>H<sub>6</sub>AsO<sub>2</sub>).





Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total 5		C 2		0	0
3	В	1	Total 5	As 1	C 2	O 2	0	0

#### • Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	23	Total O 23 23	0	0
4	В	14	Total O 14 14	0	0



## 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.



## 4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 31	Depositor
Cell constants	59.89Å 59.89Å 129.87Å	Donositon
a, b, c, $\alpha$ , $\beta$ , $\gamma$	$90.00^{\circ}$ $90.00^{\circ}$ $120.00^{\circ}$	Depositor
Resolution (Å)	33.24 - 2.60	Depositor
rtesolution (A)	33.24 - 2.60	EDS
% Data completeness	95.9 (33.24-2.60)	Depositor
(in resolution range)	95.9 (33.24-2.60)	EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$< I/\sigma(I) > 1$	5.01 (at 2.61Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
$R, R_{free}$	0.172 , $0.240$	Depositor
	0.177 , $0.242$	DCC
$R_{free}$ test set	1592 reflections (9.97%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	45.2	Xtriage
Anisotropy	0.614	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$ , $B_{sol}(Å^2)$	$0.32\;,27.7$	EDS
L-test for twinning <sup>2</sup>	$< L > = 0.45, < L^2> = 0.28$	Xtriage
	0.036 for -h,-k,l	
Estimated twinning fraction	0.166  for h,-h-k,-l	Xtriage
	0.049  for -k,-h,-l	
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	3700	wwPDB-VP
Average B, all atoms $(\mathring{A}^2)$	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 5.41% of the height of the origin peak. No significant pseudotranslation is detected.

<sup>&</sup>lt;sup>2</sup>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.



<sup>&</sup>lt;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, CAC

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		
IVIOI	Chain	RMSZ	# Z  > 5	RMSZ	# Z  > 5	
1	A	0.45	0/1887	0.62	0/2580	
1	В	0.46	0/1869	0.65	$1/2554 \ (0.0\%)$	
All	All	0.46	0/3756	0.64	1/5134 (0.0%)	

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	$\mathbf{Z}$	$\mathbf{Observed}(^{o})$	$\operatorname{Ideal}({}^{o})$
1	В	88	LEU	CA-CB-CG	6.90	131.16	115.30

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	A	1834	0	1769	21	0
1	В	1817	0	1753	23	0
2	A	1	0	0	0	0
2	В	1	0	0	0	0
3	A	5	0	0	1	0
3	В	5	0	0	0	0
4	A	23	0	0	3	0
4	В	14	0	0	1	0

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$\mathbf{Mol}$	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	3700	0	3522	43	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 6.

All (43) 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:116:ASN:HD22	1:B:116:ASN:N	1.71	0.89
1:B:116:ASN:H	1:B:119:ALA:HB2	1.40	0.86
1:B:114:ARG:HG3	1:B:122:ARG:HG2	1.63	0.80
1:A:53:ASN:OD1	1:B:53:ASN:HB3	1.85	0.75
1:A:236:LYS:O	4:A:401:HOH:O	2.06	0.71
1:A:187:VAL:O	1:A:189:TYR:N	2.22	0.69
1:A:110:GLU:O	4:A:402:HOH:O	2.14	0.65
1:A:2:THR:HG22	1:A:4:THR:H	1.62	0.64
1:B:116:ASN:N	1:B:116:ASN:ND2	2.44	0.64
1:A:120:GLU:OE2	4:A:403:HOH:O	2.16	0.60
1:B:158:ASP:HB2	1:B:159:PRO:HA	1.86	0.57
1:B:63:LEU:HD12	1:B:77:LEU:HD21	1.88	0.54
1:A:148:ILE:HG13	1:A:150:PRO:HD3	1.90	0.54
1:A:98:ASN:ND2	3:A:302:CAC:O2	2.41	0.54
1:A:72:ARG:HG3	1:A:232:ARG:HA	1.89	0.53
1:B:24:LEU:HD23	1:B:92:ALA:HB3	1.90	0.53
1:B:159:PRO:HB2	1:B:213:PRO:HG3	1.89	0.53
1:B:159:PRO:O	1:B:213:PRO:HD3	2.11	0.50
1:A:45:LEU:HD12	1:A:76:VAL:HB	1.94	0.50
1:B:160:LEU:O	1:B:210:GLU:HB3	2.12	0.50
1:A:10:ARG:O	1:A:64:ARG:NH1	2.44	0.49
1:B:111:THR:HG23	1:B:126:LEU:HB2	1.95	0.49
1:B:131:LYS:HB2	1:B:132:PRO:HD2	1.96	0.47
1:B:131:LYS:O	1:B:134:SER:HB3	2.15	0.46
1:B:169:SER:O	1:B:173:GLU:HG2	2.15	0.46
1:A:48:THR:HA	1:A:79:VAL:HB	1.97	0.45
1:A:52:GLU:HG2	1:A:158:ASP:HB2	1.99	0.44
1:A:14:PRO:HA	1:A:15:PRO:HD3	1.92	0.43
1:A:50:GLY:O	1:A:84:CYS:HA	2.19	0.43
1:B:52:GLU:HG2	1:B:158:ASP:OD2	2.19	0.43
1:A:171:LEU:HD13	1:A:233:TRP:CE2	2.54	0.43
1:B:208:THR:HG22	1:B:210:GLU:HG2	2.01	0.43
1:A:134:SER:OG	1:A:135:GLU:OE2	2.34	0.42

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Atom-1	Atom-2	$egin{aligned} &  ext{Interatomic} \ &  ext{distance} \ &  ext{(Å)} \end{aligned}$	Clash overlap (Å)
1:A:171:LEU:HD13	1:A:233:TRP:CD2	2.55	0.41
1:B:1:MET:N	4:B:401:HOH:O	2.30	0.41
1:A:72:ARG:HD2	1:A:232:ARG:O	2.20	0.41
1:B:48:THR:HA	1:B:79:VAL:HB	2.01	0.41
1:B:72:ARG:NH1	1:B:234:HIS:HB3	2.35	0.41
1:B:233:TRP:CZ3	1:B:235:PRO:HD3	2.55	0.41
1:A:34:ALA:HA	1:A:35:PRO:HD3	1.94	0.41
1:B:116:ASN:HD22	1:B:116:ASN:H	1.60	0.41
1:A:153:VAL:HG11	1:A:195:PHE:CZ	2.55	0.41
1:B:135:GLU:O	1:B:139:GLN:HG3	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	235/237~(99%)	220 (94%)	12 (5%)	3 (1%)	12	24	
1	В	$230/237 \ (97\%)$	216 (94%)	13 (6%)	1 (0%)	34	57	
All	All	465/474 (98%)	436 (94%)	25 (5%)	4 (1%)	17	35	

#### All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	188	GLY
1	В	168	HIS
1	A	102	PRO
1	A	159	PRO



#### 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	196/196 (100%)	189 (96%)	7 (4%)	35	61	
1	В	194/196 (99%)	189 (97%)	5 (3%)	46	72	
All	All	390/392 (100%)	378 (97%)	12 (3%)	40	66	

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	55	SER
1	A	84	CYS
1	A	85	$\operatorname{GLN}$
1	A	121	GLU
1	A	184	VAL
1	A	185	THR
1	В	40	GLU
1	В	67	THR
1	В	116	ASN
1	В	191	THR
1	В	220	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	В	116	ASN

#### 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)

Of 4 ligands modelled in this entry, 2 are monoatomic - leaving 2 for Mogul analysis.

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		$\mathbf{n} \mid \mathbf{Res} \mid \mathbf{Link} \mid$		B	Bond lengths			ond ang	gles	
MIOI	Mol Type Chain	Res Lin	LIIIK	Counts	RMSZ	# Z >2	Counts	RMSZ	# Z >2	
3	CAC	A	302	2	0,4,4	-	-	0,6,6	-	-
3	CAC	В	302	2	0,4,4	-	-	0,6,6	-	-

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.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	302	CAC	1	0

#### 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	<RSRZ $>$	$\#\mathrm{RSRZ}{>}2$		$OWAB(A^2)$	Q<0.9	
1	A	237/237 (100%)	-0.26	2 (0%)	86	84	36, 51, 68, 97	0
1	В	$234/237 \ (98\%)$	-0.32	3 (1%)	77	73	37, 49, 75, 103	0
All	All	471/474 (99%)	-0.29	5 (1%)	80	78	36, 50, 70, 103	0

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	В	236	LYS	3.5
1	В	121	GLU	2.8
1	A	3	VAL	2.3
1	A	237	ASP	2.2
1	В	189	TYR	2.1

#### 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	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{\mathring{A}}^2)$	Q<0.9
3	CAC	A	302	5/5	0.73	0.17	52,55,60,105	5
3	CAC	В	302	5/5	0.80	0.21	48,59,69,97	5
2	ZN	В	301	1/1	0.90	0.15	75,75,75,75	1
2	ZN	A	301	1/1	0.97	0.09	72,72,72,72	1

# 6.5 Other polymers (i)

There are no such residues in this entry.

