



# wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 1, 2016 – 05:48 PM GMT

PDB ID : 4JHM  
Title : Crystal structure of a putative mandelate racemase/muconate lactonizing enzyme from *Pseudovibrio* sp.  
Authors : Hegde, R.P.; Toro, R.; Burley, S.K.; Almo, S.C.; Ramagopal, U.A.; New York Structural Genomics Research Consortium (NYSGRG)  
Deposited on : 2013-03-05  
Resolution : 2.80 Å(reported)

This is a wwPDB X-ray Structure 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/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

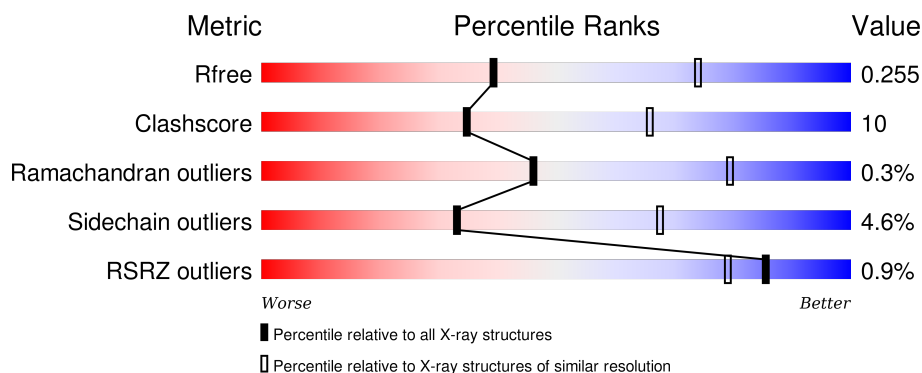
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.80 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	2393 (2.80-2.80)
Clashscore	102246	2827 (2.80-2.80)
Ramachandran outliers	100387	2782 (2.80-2.80)
Sidechain outliers	100360	2784 (2.80-2.80)
RSRZ outliers	91569	2404 (2.80-2.80)

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	384	<div> <div>76%</div> <div>20%</div> <div>• •</div> </div>
1	B	384	<div> <div>81%</div> <div>15%</div> <div>• •</div> </div>
1	C	384	<div> <div>%</div> <div>75%</div> <div>20%</div> <div>• •</div> </div>
1	D	384	<div> <div>%</div> <div>74%</div> <div>21%</div> <div>• •</div> </div>
1	E	384	<div> <div>%</div> <div>73%</div> <div>22%</div> <div>• •</div> </div>

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Mol	Chain	Length	Quality of chain
1	F	384	<div><div></div><div>73%22%<div><div></div><div></div></div></div></div>
1	G	384	<div>%<div><div></div><div>72%23%<div><div></div><div></div></div></div></div></div>
1	H	384	<div>3%<div><div></div><div>76%19%<div><div></div><div></div></div></div></div></div>

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 23568 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 Mandelate racemase / muconate lactonizing enzyme, C-terminal domain protein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	375	Total	C	N	O	S	Se	0	0	0
			2946	1876	482	566	10	12			
1	B	375	Total	C	N	O	S	Se	0	1	0
			2960	1886	484	568	10	12			
1	C	374	Total	C	N	O	S	Se	0	0	0
			2940	1874	479	565	10	12			
1	D	375	Total	C	N	O	S	Se	0	0	0
			2942	1873	481	566	10	12			
1	E	374	Total	C	N	O	S	Se	0	2	0
			2951	1880	480	569	10	12			
1	F	375	Total	C	N	O	S	Se	0	1	0
			2945	1875	479	569	10	12			
1	G	375	Total	C	N	O	S	Se	0	0	0
			2944	1876	480	566	10	12			
1	H	374	Total	C	N	O	S	Se	0	0	0
			2932	1868	477	565	10	12			

There are 88 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	MSE	-	EXPRESSION TAG	UNP B6R2Z8
A	0	SER	-	EXPRESSION TAG	UNP B6R2Z8
A	1	LEU	-	EXPRESSION TAG	UNP B6R2Z8
A	375	GLU	-	EXPRESSION TAG	UNP B6R2Z8
A	376	GLY	-	EXPRESSION TAG	UNP B6R2Z8
A	377	HIS	-	EXPRESSION TAG	UNP B6R2Z8
A	378	HIS	-	EXPRESSION TAG	UNP B6R2Z8
A	379	HIS	-	EXPRESSION TAG	UNP B6R2Z8
A	380	HIS	-	EXPRESSION TAG	UNP B6R2Z8
A	381	HIS	-	EXPRESSION TAG	UNP B6R2Z8
A	382	HIS	-	EXPRESSION TAG	UNP B6R2Z8
B	-1	MSE	-	EXPRESSION TAG	UNP B6R2Z8

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Chain	Residue	Modelled	Actual	Comment	Reference
B	0	SER	-	EXPRESSION TAG	UNP B6R2Z8
B	1	LEU	-	EXPRESSION TAG	UNP B6R2Z8
B	375	GLU	-	EXPRESSION TAG	UNP B6R2Z8
B	376	GLY	-	EXPRESSION TAG	UNP B6R2Z8
B	377	HIS	-	EXPRESSION TAG	UNP B6R2Z8
B	378	HIS	-	EXPRESSION TAG	UNP B6R2Z8
B	379	HIS	-	EXPRESSION TAG	UNP B6R2Z8
B	380	HIS	-	EXPRESSION TAG	UNP B6R2Z8
B	381	HIS	-	EXPRESSION TAG	UNP B6R2Z8
B	382	HIS	-	EXPRESSION TAG	UNP B6R2Z8
C	-1	MSE	-	EXPRESSION TAG	UNP B6R2Z8
C	0	SER	-	EXPRESSION TAG	UNP B6R2Z8
C	1	LEU	-	EXPRESSION TAG	UNP B6R2Z8
C	375	GLU	-	EXPRESSION TAG	UNP B6R2Z8
C	376	GLY	-	EXPRESSION TAG	UNP B6R2Z8
C	377	HIS	-	EXPRESSION TAG	UNP B6R2Z8
C	378	HIS	-	EXPRESSION TAG	UNP B6R2Z8
C	379	HIS	-	EXPRESSION TAG	UNP B6R2Z8
C	380	HIS	-	EXPRESSION TAG	UNP B6R2Z8
C	381	HIS	-	EXPRESSION TAG	UNP B6R2Z8
C	382	HIS	-	EXPRESSION TAG	UNP B6R2Z8
D	-1	MSE	-	EXPRESSION TAG	UNP B6R2Z8
D	0	SER	-	EXPRESSION TAG	UNP B6R2Z8
D	1	LEU	-	EXPRESSION TAG	UNP B6R2Z8
D	375	GLU	-	EXPRESSION TAG	UNP B6R2Z8
D	376	GLY	-	EXPRESSION TAG	UNP B6R2Z8
D	377	HIS	-	EXPRESSION TAG	UNP B6R2Z8
D	378	HIS	-	EXPRESSION TAG	UNP B6R2Z8
D	379	HIS	-	EXPRESSION TAG	UNP B6R2Z8
D	380	HIS	-	EXPRESSION TAG	UNP B6R2Z8
D	381	HIS	-	EXPRESSION TAG	UNP B6R2Z8
D	382	HIS	-	EXPRESSION TAG	UNP B6R2Z8
E	-1	MSE	-	EXPRESSION TAG	UNP B6R2Z8
E	0	SER	-	EXPRESSION TAG	UNP B6R2Z8
E	1	LEU	-	EXPRESSION TAG	UNP B6R2Z8
E	375	GLU	-	EXPRESSION TAG	UNP B6R2Z8
E	376	GLY	-	EXPRESSION TAG	UNP B6R2Z8
E	377	HIS	-	EXPRESSION TAG	UNP B6R2Z8
E	378	HIS	-	EXPRESSION TAG	UNP B6R2Z8
E	379	HIS	-	EXPRESSION TAG	UNP B6R2Z8
E	380	HIS	-	EXPRESSION TAG	UNP B6R2Z8
E	381	HIS	-	EXPRESSION TAG	UNP B6R2Z8

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Chain	Residue	Modelled	Actual	Comment	Reference
E	382	HIS	-	EXPRESSION TAG	UNP B6R2Z8
F	-1	MSE	-	EXPRESSION TAG	UNP B6R2Z8
F	0	SER	-	EXPRESSION TAG	UNP B6R2Z8
F	1	LEU	-	EXPRESSION TAG	UNP B6R2Z8
F	375	GLU	-	EXPRESSION TAG	UNP B6R2Z8
F	376	GLY	-	EXPRESSION TAG	UNP B6R2Z8
F	377	HIS	-	EXPRESSION TAG	UNP B6R2Z8
F	378	HIS	-	EXPRESSION TAG	UNP B6R2Z8
F	379	HIS	-	EXPRESSION TAG	UNP B6R2Z8
F	380	HIS	-	EXPRESSION TAG	UNP B6R2Z8
F	381	HIS	-	EXPRESSION TAG	UNP B6R2Z8
F	382	HIS	-	EXPRESSION TAG	UNP B6R2Z8
G	-1	MSE	-	EXPRESSION TAG	UNP B6R2Z8
G	0	SER	-	EXPRESSION TAG	UNP B6R2Z8
G	1	LEU	-	EXPRESSION TAG	UNP B6R2Z8
G	375	GLU	-	EXPRESSION TAG	UNP B6R2Z8
G	376	GLY	-	EXPRESSION TAG	UNP B6R2Z8
G	377	HIS	-	EXPRESSION TAG	UNP B6R2Z8
G	378	HIS	-	EXPRESSION TAG	UNP B6R2Z8
G	379	HIS	-	EXPRESSION TAG	UNP B6R2Z8
G	380	HIS	-	EXPRESSION TAG	UNP B6R2Z8
G	381	HIS	-	EXPRESSION TAG	UNP B6R2Z8
G	382	HIS	-	EXPRESSION TAG	UNP B6R2Z8
H	-1	MSE	-	EXPRESSION TAG	UNP B6R2Z8
H	0	SER	-	EXPRESSION TAG	UNP B6R2Z8
H	1	LEU	-	EXPRESSION TAG	UNP B6R2Z8
H	375	GLU	-	EXPRESSION TAG	UNP B6R2Z8
H	376	GLY	-	EXPRESSION TAG	UNP B6R2Z8
H	377	HIS	-	EXPRESSION TAG	UNP B6R2Z8
H	378	HIS	-	EXPRESSION TAG	UNP B6R2Z8
H	379	HIS	-	EXPRESSION TAG	UNP B6R2Z8
H	380	HIS	-	EXPRESSION TAG	UNP B6R2Z8
H	381	HIS	-	EXPRESSION TAG	UNP B6R2Z8
H	382	HIS	-	EXPRESSION TAG	UNP B6R2Z8

- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	1	Total O 1 1	0	0
2	C	2	Total O 2 2	0	0

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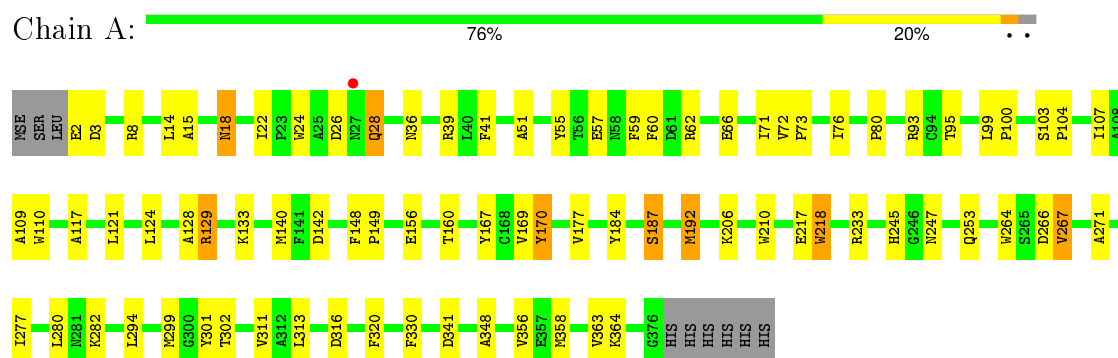
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	D	1	Total 1	O 1	0	0
2	E	1	Total 1	O 1	0	0
2	F	1	Total 1	O 1	0	0
2	G	2	Total 2	O 2	0	0

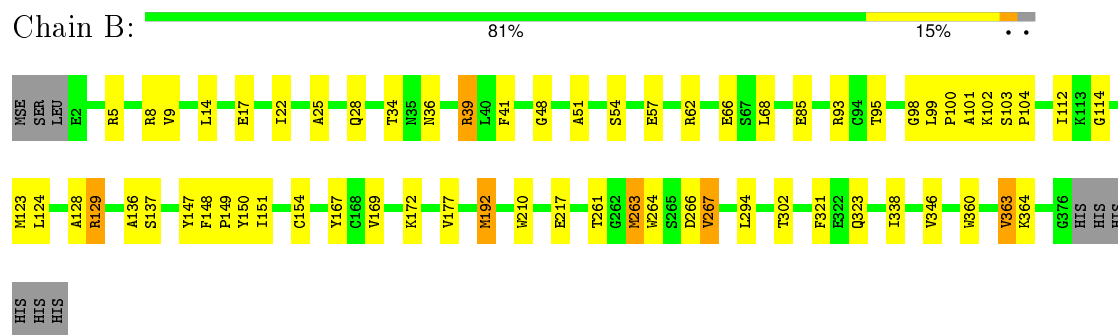
### 3 Residue-property plots

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.

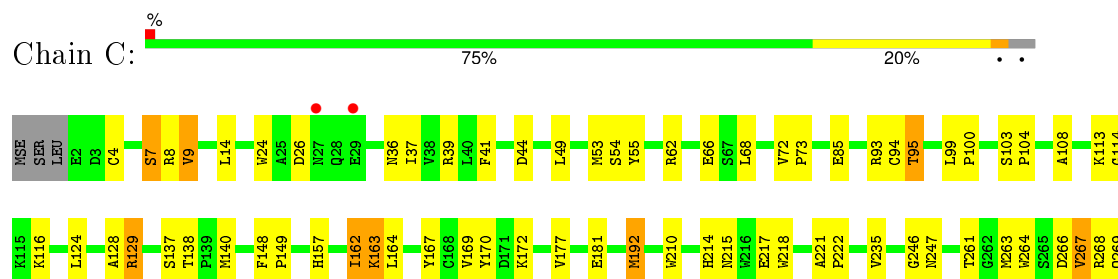
- Molecule 1: Mandelate racemase / muconate lactonizing enzyme, C-terminal domain protein



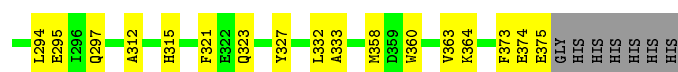
- Molecule 1: Mandelate racemase / muconate lactonizing enzyme, C-terminal domain protein



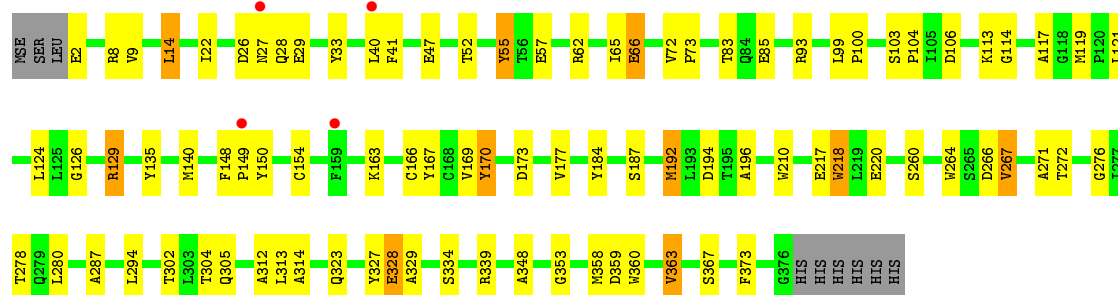
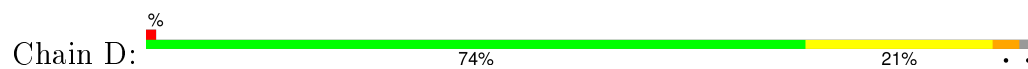
- Molecule 1: Mandelate racemase / muconate lactonizing enzyme, C-terminal domain protein



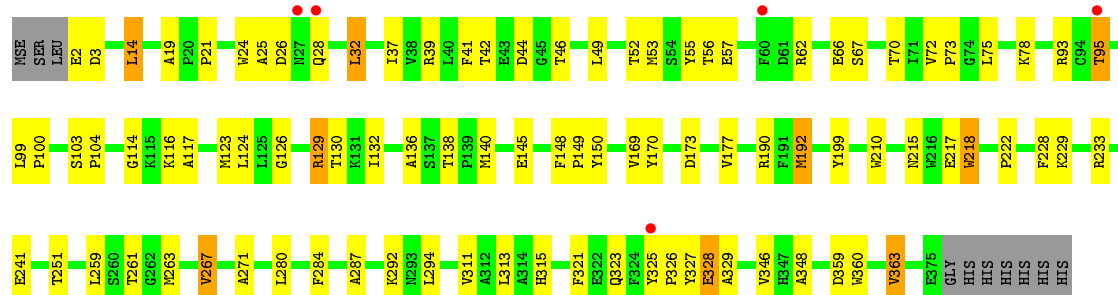




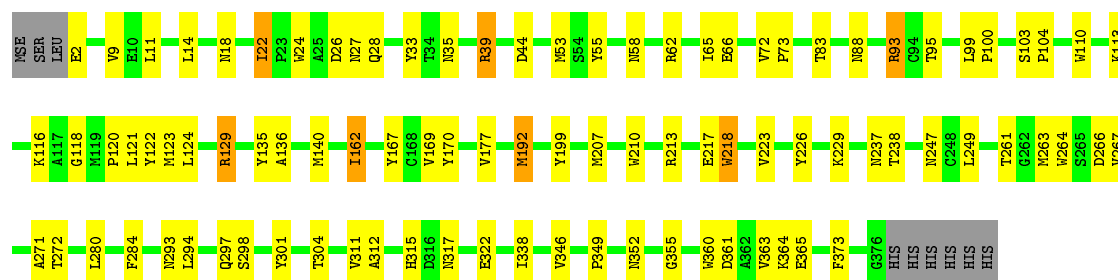
- Molecule 1: Mandelate racemase / muconate lactonizing enzyme, C-terminal domain protein



- Molecule 1: Mandelate racemase / muconate lactonizing enzyme, C-terminal domain protein

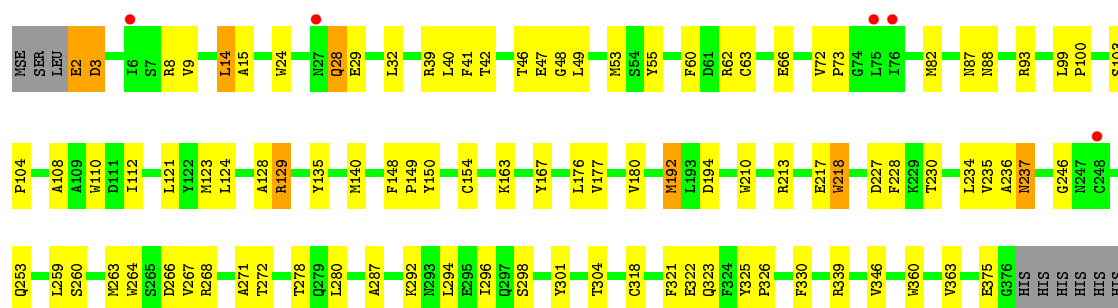


- Molecule 1: Mandelate racemase / muconate lactonizing enzyme, C-terminal domain protein

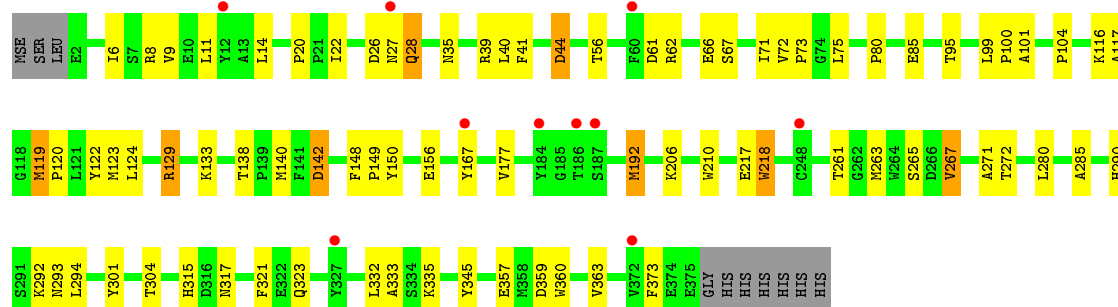
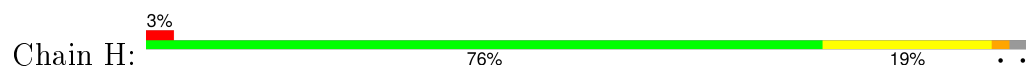


- Molecule 1: Mandelate racemase / muconate lactonizing enzyme, C-terminal domain protein





- Molecule 1: Mandelate racemase / muconate lactonizing enzyme, C-terminal domain protein



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	201.79Å 117.31Å 143.13Å 90.00° 99.77° 90.00°	Depositor
Resolution (Å)	39.41 – 2.80 39.41 – 2.80	Depositor EDS
% Data completeness (in resolution range)	98.6 (39.41-2.80) 98.6 (39.41-2.80)	Depositor EDS
$R_{merge}$	0.05	Depositor
$R_{sym}$	0.10	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.66 (at 2.81Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, $R_{free}$	0.194 , 0.253 0.195 , 0.255	Depositor DCC
$R_{free}$ test set	3989 reflections (5.25%)	DCC
Wilson B-factor (Å <sup>2</sup> )	66.8	Xtriage
Anisotropy	0.052	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.28 , 29.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.52$ , $\langle L^2 \rangle = 0.35$	Xtriage
Outliers	0 of 79901 reflections	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	23568	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	85.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.55% of the height of the origin peak. No significant pseudotranslation is detected.*

<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 [i](#)

### 5.1 Standard geometry [i](#)

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	0.27	0/3009	0.46	0/4072
1	B	0.28	0/3026	0.49	0/4092
1	C	0.28	0/3003	0.46	0/4064
1	D	0.26	0/3005	0.44	0/4068
1	E	0.26	0/3017	0.44	0/4084
1	F	0.26	0/3008	0.45	0/4073
1	G	0.27	0/3007	0.45	0/4069
1	H	0.26	0/2995	0.44	0/4056
All	All	0.27	0/24070	0.46	0/32578

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	A	2946	0	2808	57	0
1	B	2960	0	2835	48	0
1	C	2940	0	2805	60	0
1	D	2942	0	2797	63	0
1	E	2951	0	2806	70	0
1	F	2945	0	2790	63	0
1	G	2944	0	2808	67	0
1	H	2932	0	2783	65	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	1	0	0	0	0
2	C	2	0	0	0	0
2	D	1	0	0	0	0
2	E	1	0	0	0	0
2	F	1	0	0	0	0
2	G	2	0	0	0	0
All	All	23568	0	22432	457	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 10.

The worst 5 of 457 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:25:ALA:HB3	1:B:28:GLN:HG2	1.58	0.85
1:C:177:VAL:HG12	1:C:210:TRP:CH2	2.12	0.84
1:E:321:PHE:HE2	1:E:323:GLN:HG2	1.46	0.81
1:E:66:GLU:HG2	1:F:62:ARG:HB2	1.61	0.81
1:D:272:THR:HG21	1:D:304:THR:HG23	1.63	0.81

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	373/384 (97%)	353 (95%)	18 (5%)	2 (0%)	34	69
1	B	373/384 (97%)	359 (96%)	14 (4%)	0	100	100
1	C	372/384 (97%)	351 (94%)	19 (5%)	2 (0%)	34	69
1	D	373/384 (97%)	348 (93%)	22 (6%)	3 (1%)	24	58
1	E	373/384 (97%)	353 (95%)	18 (5%)	2 (0%)	34	69

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	F	373/384 (97%)	352 (94%)	21 (6%)	0	100	100
1	G	373/384 (97%)	355 (95%)	17 (5%)	1 (0%)	46	79
1	H	372/384 (97%)	357 (96%)	15 (4%)	0	100	100
All	All	2982/3072 (97%)	2828 (95%)	144 (5%)	10 (0%)	46	79

5 of 10 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	328	GLU
1	D	170	TYR
1	E	328	GLU
1	C	95	THR
1	D	196	ALA

### 5.3.2 Protein sidechains ⓘ

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	309/307 (101%)	291 (94%)	18 (6%)	25	57
1	B	312/307 (102%)	298 (96%)	14 (4%)	34	68
1	C	309/307 (101%)	295 (96%)	14 (4%)	34	68
1	D	308/307 (100%)	296 (96%)	12 (4%)	39	74
1	E	310/307 (101%)	297 (96%)	13 (4%)	36	71
1	F	308/307 (100%)	293 (95%)	15 (5%)	31	65
1	G	309/307 (101%)	293 (95%)	16 (5%)	29	62
1	H	307/307 (100%)	294 (96%)	13 (4%)	36	71
All	All	2472/2456 (101%)	2357 (95%)	115 (5%)	33	67

5 of 115 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	119	MSE

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Mol	Chain	Res	Type
1	E	95	THR
1	H	85	GLU
1	D	129	ARG
1	E	2[A]	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 33 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	165	HIS
1	D	323	GLN
1	G	323	GLN
1	C	247	ASN
1	C	290	HIS

### 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](#)

There are no ligands in this entry.

## 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 ⓘ

### 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	363/384 (94%)	-0.35	1 (0%) 94 92	50, 78, 108, 151	0
1	B	363/384 (94%)	-0.42	0 100 100	47, 68, 101, 145	0
1	C	362/384 (94%)	-0.40	2 (0%) 90 86	48, 74, 109, 159	0
1	D	363/384 (94%)	-0.16	4 (1%) 82 74	53, 89, 122, 155	0
1	E	362/384 (94%)	-0.12	5 (1%) 78 69	57, 92, 116, 152	0
1	F	363/384 (94%)	-0.35	0 100 100	56, 87, 115, 146	0
1	G	363/384 (94%)	-0.15	5 (1%) 78 69	56, 87, 113, 146	0
1	H	362/384 (94%)	-0.11	10 (2%) 56 44	58, 92, 118, 155	0
All	All	2901/3072 (94%)	-0.26	27 (0%) 85 79	47, 84, 116, 159	0

The worst 5 of 27 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	27	ASN	5.1
1	A	27	ASN	3.9
1	H	186	THR	3.5
1	E	28[A]	GLN	3.2
1	H	248	CYS	3.1

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

There are no non-standard protein/DNA/RNA residues in this entry.

### 6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.



## 6.4 Ligands

There are no ligands in this entry.

## 6.5 Other polymers

There are no such residues in this entry.