



# wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 9, 2018 – 11:38 am GMT

PDB ID : 2D4A  
Title : Structure of the malate dehydrogenase from *Aeropyrum pernix*  
Authors : Kawakami, R.; Sakuraba, H.; Tsuge, H.; Ohshima, T.  
Deposited on : 2005-10-12  
Resolution : 2.87 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

<https://www.wwpdb.org/validation/2017/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
Xtriage (Phenix)	:	1.13
EDS	:	trunk30967
Percentile statistics	:	20171227.v01 (using entries in the PDB archive December 27th 2017)
Refmac	:	5.8.0158
CCP4	:	7.0 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk30967

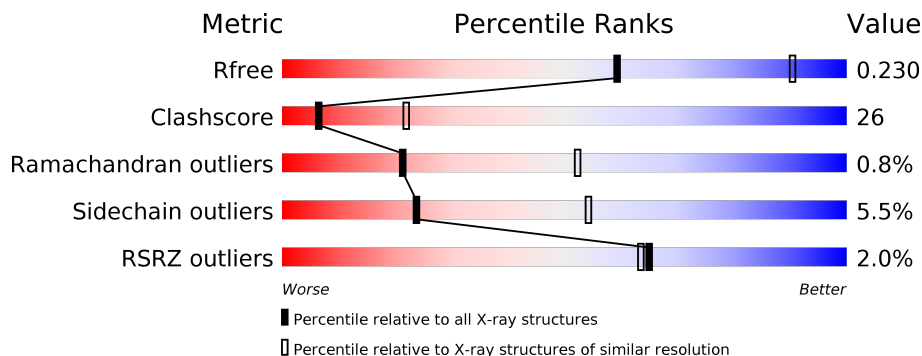
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.87 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))
$R_{free}$	111664	2330 (2.90-2.86)
Clashscore	122126	2579 (2.90-2.86)
Ramachandran outliers	120053	2524 (2.90-2.86)
Sidechain outliers	120020	2527 (2.90-2.86)
RSRZ outliers	108989	2272 (2.90-2.86)

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	308	<div> <div>0.5%</div> <div>56%</div> <div>35%</div> <div>5%</div> <div>• •</div> </div>
1	B	308	<div> <div>2%</div> <div>61%</div> <div>30%</div> <div>7%</div> <div>•</div> </div>
1	C	308	<div> <div>2%</div> <div>50%</div> <div>39%</div> <div>7%</div> <div>• •</div> </div>
1	D	308	<div> <div>3%</div> <div>59%</div> <div>36%</div> <div>5%</div> <div>•</div> </div>

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 9292 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 Malate dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	301	Total	C	N	O	S	0	0	0
			2289	1455	383	438	13			
1	D	308	Total	C	N	O	S	0	0	0
			2346	1489	397	446	14			
1	C	299	Total	C	N	O	S	0	0	0
			2274	1446	384	431	13			
1	A	298	Total	C	N	O	S	0	0	0
			2262	1440	379	430	13			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	INITIATING METHIONINE	UNP Q9YEA1
B	1	MET	-	INITIATING METHIONINE	UNP Q9YEA1
C	1	MET	-	INITIATING METHIONINE	UNP Q9YEA1
D	1	MET	-	INITIATING METHIONINE	UNP Q9YEA1

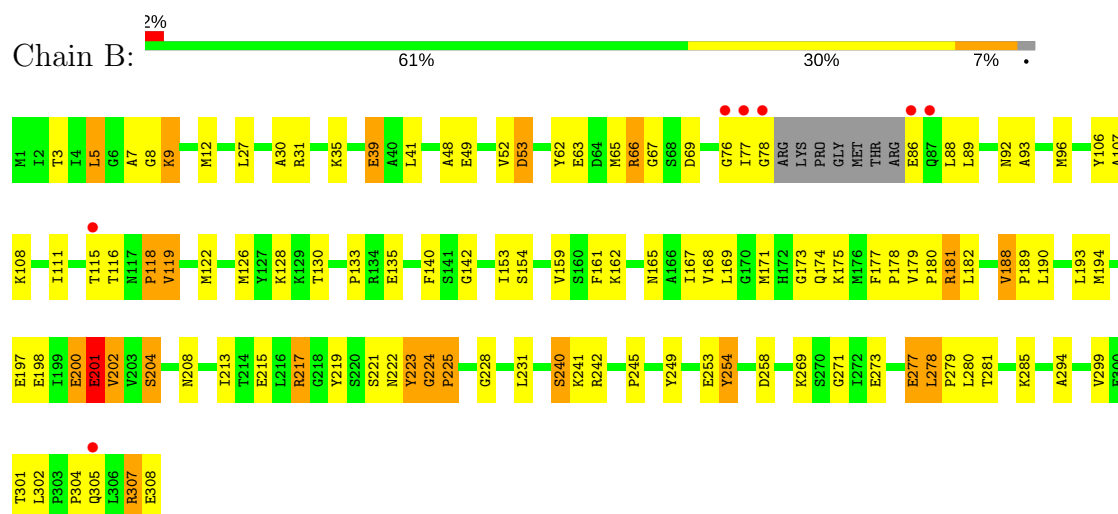
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	42	Total	O	0	0
			42	42		
2	D	34	Total	O	0	0
			34	34		
2	C	17	Total	O	0	0
			17	17		
2	A	28	Total	O	0	0
			28	28		

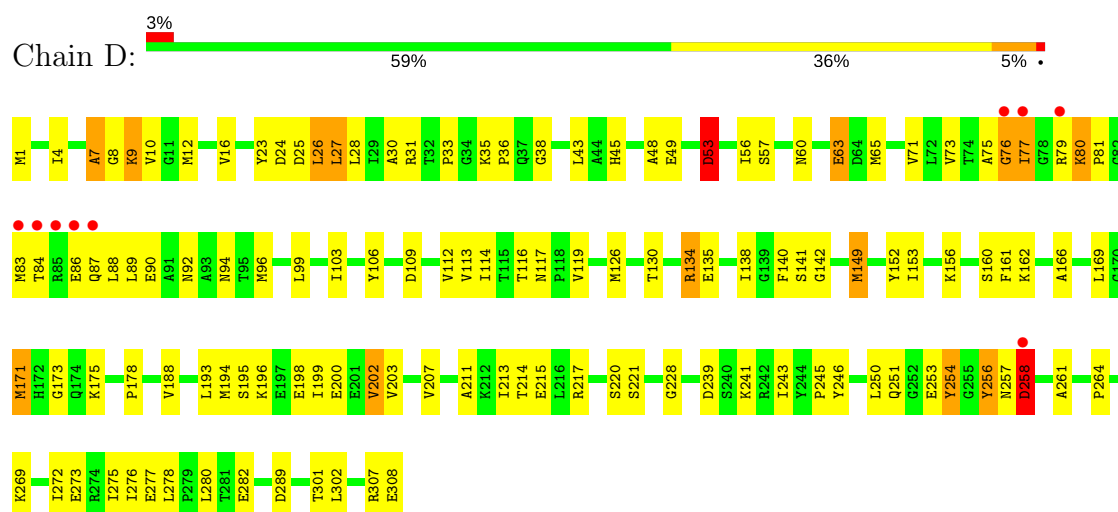
### 3 Residue-property plots

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

#### • Molecule 1: Malate dehydrogenase

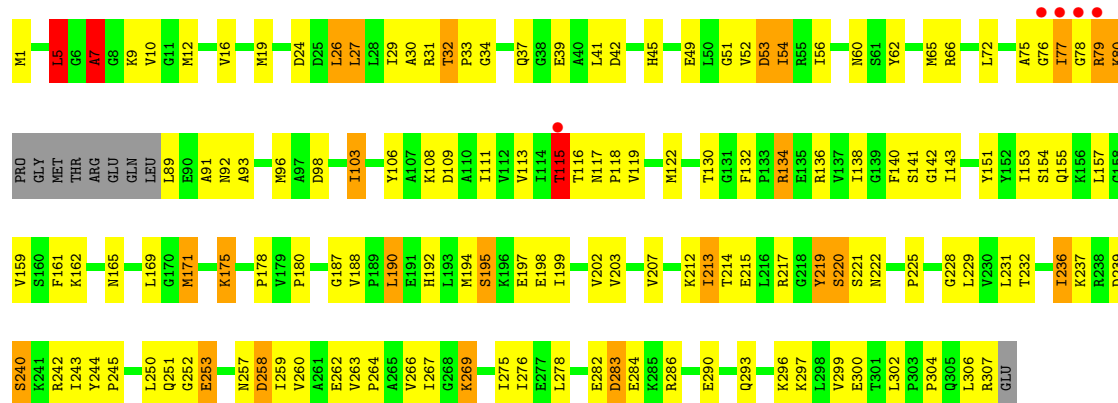


#### • Molecule 1: Malate dehydrogenase

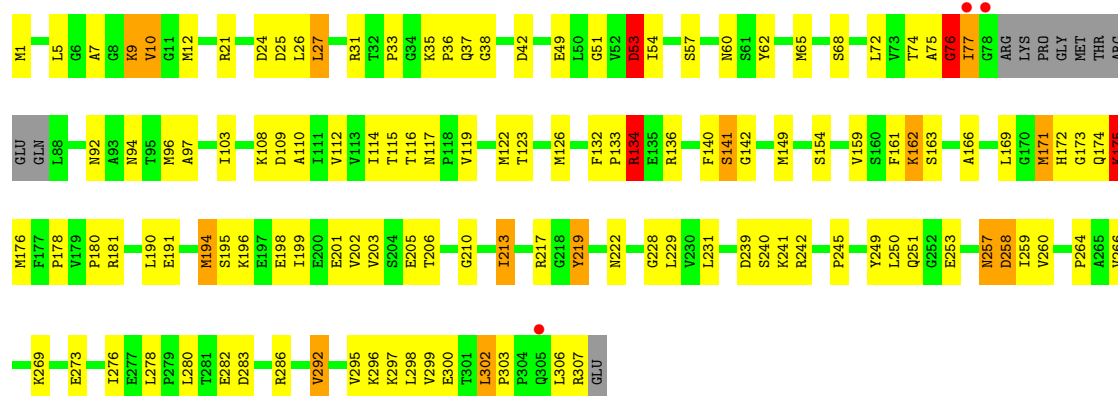


#### • Molecule 1: Malate dehydrogenase





● Molecule 1: Malate dehydrogenase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	78.17Å 84.01Å 216.15Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.92 – 2.87 5.15 – 2.87	Depositor EDS
% Data completeness (in resolution range)	(Not available) (19.92-2.87) 99.8 (5.15-2.87)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.86 (at 2.89Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.205 , 0.240 0.195 , 0.230	Depositor DCC
$R_{free}$ test set	2707 reflections (9.95%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	41.8	Xtriage
Anisotropy	0.086	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.86 , 113.5	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	9292	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 11.49% 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, 0.333 respectively for untwinned datasets, and 0.375, 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	0.65	7/2295 (0.3%)	1.08	20/3100 (0.6%)
1	B	0.88	13/2322 (0.6%)	1.31	37/3136 (1.2%)
1	C	0.70	10/2307 (0.4%)	1.04	16/3114 (0.5%)
1	D	0.63	4/2381 (0.2%)	0.97	16/3215 (0.5%)
All	All	0.72	34/9305 (0.4%)	1.10	89/12565 (0.7%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	3
1	B	0	3
1	C	0	4
1	D	0	2
All	All	0	12

The worst 5 of 34 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	254	TYR	C-N	8.37	1.48	1.33
1	C	5	LEU	C-N	7.81	1.47	1.33
1	C	253	GLU	C-N	-7.72	1.16	1.34
1	A	134	ARG	C-O	-7.71	1.08	1.23
1	B	217	ARG	C-N	-7.65	1.19	1.33

The worst 5 of 89 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	7	ALA	O-C-N	-12.69	101.64	123.20
1	C	115	THR	O-C-N	-12.28	103.06	122.70
1	C	219	TYR	O-C-N	-11.90	103.66	122.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	201	GLU	O-C-N	-11.79	103.84	122.70
1	C	240	SER	O-C-N	-11.14	104.88	122.70

There are no chirality outliers.

5 of 12 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	201	GLU	Mainchain
1	B	202	VAL	Mainchain
1	B	278	LEU	Mainchain
1	D	258	ASP	Mainchain
1	D	8	GLY	Mainchain

## 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	2262	0	2334	111	1
1	B	2289	0	2353	118	2
1	C	2274	0	2349	146	0
1	D	2346	0	2422	132	1
2	A	28	0	0	1	0
2	B	42	0	0	1	0
2	C	17	0	0	0	0
2	D	34	0	0	1	0
All	All	9292	0	9458	475	3

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 26.

The worst 5 of 475 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:174:GLN:NE2	1:B:175:LYS:HZ2	1.34	1.24
1:B:174:GLN:NE2	1:B:175:LYS:NZ	1.93	1.17

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:76:GLY:HA3	1:B:116:THR:HG23	1.32	1.09
1:A:76:GLY:HA3	1:A:116:THR:HG23	1.33	1.06
1:A:259:ILE:HD13	1:A:292:VAL:HG13	1.44	1.00

All (3) 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	Interatomic distance (Å)	Clash overlap (Å)
1:B:108:LYS:NZ	1:B:200:GLU:OE1[4_557]	1.86	0.34
1:B:133:PRO:CA	1:B:215:GLU:OE1[4_557]	1.99	0.21
1:D:87:GLN:NE2	1:A:181:ARG:NH1[3_756]	2.02	0.18

## 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	294/308 (96%)	276 (94%)	15 (5%)	3 (1%)	17	46
1	B	297/308 (96%)	280 (94%)	16 (5%)	1 (0%)	43	73
1	C	295/308 (96%)	279 (95%)	13 (4%)	3 (1%)	17	46
1	D	306/308 (99%)	287 (94%)	17 (6%)	2 (1%)	24	56
All	All	1192/1232 (97%)	1122 (94%)	61 (5%)	9 (1%)	21	52

5 of 9 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	53	ASP
1	C	220	SER
1	B	181	ARG
1	C	53	ASP
1	C	75	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	241/250 (96%)	229 (95%)	12 (5%)	27	58
1	B	244/250 (98%)	234 (96%)	10 (4%)	33	66
1	C	242/250 (97%)	224 (93%)	18 (7%)	15	39
1	D	250/250 (100%)	236 (94%)	14 (6%)	23	53
All	All	977/1000 (98%)	923 (94%)	54 (6%)	24	54

5 of 54 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	5	LEU
1	C	80	LYS
1	A	175	LYS
1	C	26	LEU
1	C	31	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 17 such sidechains are listed below:

Mol	Chain	Res	Type
1	D	174	GLN
1	D	208	ASN
1	C	165	ASN
1	D	172	HIS
1	C	251	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](#)

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](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	B	3
1	C	2
1	A	1
1	D	1

The worst 5 of 7 chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	C	219:TYR	C	220:SER	N	1.20
1	A	123:THR	C	124:TYR	N	1.20
1	B	217:ARG	C	218:GLY	N	1.19
1	B	52:VAL	C	53:ASP	N	1.18
1	B	240:SER	C	241:LYS	N	1.18

## 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<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	298/308 (96%)	-0.57	3 (1%) 82 81	21, 35, 58, 74	0
1	B	301/308 (97%)	-0.44	7 (2%) 60 58	18, 34, 53, 79	0
1	C	299/308 (97%)	-0.49	5 (1%) 70 69	23, 39, 62, 82	0
1	D	308/308 (100%)	-0.61	9 (2%) 51 47	20, 34, 59, 71	0
All	All	1206/1232 (97%)	-0.53	24 (1%) 65 63	18, 35, 59, 82	0

The worst 5 of 24 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	87	GLN	5.7
1	D	85	ARG	4.2
1	D	84	THR	4.1
1	D	87	GLN	3.2
1	B	305	GLN	3.2

### 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](#)

There are no ligands in this entry.

## 6.5 Other polymers [i](#)

There are no such residues in this entry.