



Full wwPDB X-ray Structure Validation Report ⓘ

May 23, 2020 – 04:17 pm BST

PDB ID : 3QBP
Title : Crystal structure of fumarase Fum from Mycobacterium marinum
Authors : Seattle Structural Genomics Center for Infectious Disease (SSGCID)
Deposited on : 2011-01-13
Resolution : 1.85 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 ⓘ symbol.

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	:	2.11
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)	:	2.11

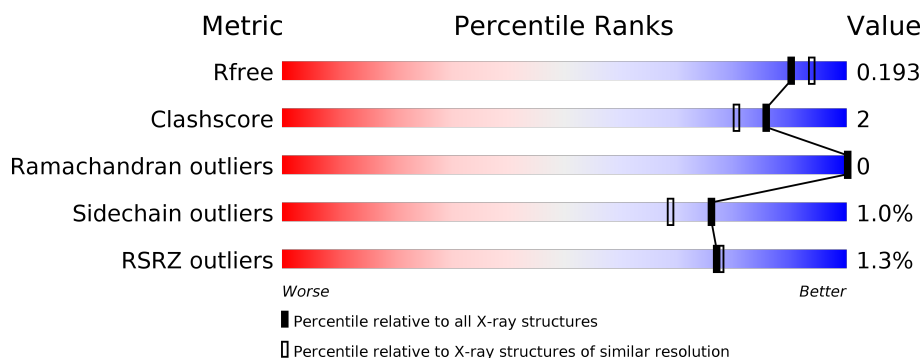
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.85 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}	130704	2469 (1.86-1.86)
Clashscore	141614	2625 (1.86-1.86)
Ramachandran outliers	138981	2592 (1.86-1.86)
Sidechain outliers	138945	2592 (1.86-1.86)
RSRZ outliers	127900	2436 (1.86-1.86)

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 ≥ 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 $\leq 5\%$. 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	478	<div> <div>%</div> <div> <div></div> <div>89%</div> <div>8%</div> </div> </div>
1	B	478	<div> <div>%</div> <div> <div></div> <div>89%</div> <div>6%</div> </div> </div>
1	C	478	<div> <div>%</div> <div> <div></div> <div>87%</div> <div>9%</div> </div> </div>
1	D	478	<div> <div>%</div> <div> <div></div> <div>88%</div> <div>9%</div> </div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 14756 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 Fumarase Fum.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	441	Total	C	N	O	S	0	11	0
			3269	2050	575	632	12			
1	B	447	Total	C	N	O	S	0	13	0
			3315	2081	586	636	12			
1	C	437	Total	C	N	O	S	0	17	0
			3282	2058	581	632	11			
1	D	437	Total	C	N	O	S	0	6	0
			3206	2012	563	621	10			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-3	GLY	-	EXPRESSION TAG	UNP B2HT54
A	-2	PRO	-	EXPRESSION TAG	UNP B2HT54
A	-1	GLY	-	EXPRESSION TAG	UNP B2HT54
A	0	SER	-	EXPRESSION TAG	UNP B2HT54
B	-3	GLY	-	EXPRESSION TAG	UNP B2HT54
B	-2	PRO	-	EXPRESSION TAG	UNP B2HT54
B	-1	GLY	-	EXPRESSION TAG	UNP B2HT54
B	0	SER	-	EXPRESSION TAG	UNP B2HT54
C	-3	GLY	-	EXPRESSION TAG	UNP B2HT54
C	-2	PRO	-	EXPRESSION TAG	UNP B2HT54
C	-1	GLY	-	EXPRESSION TAG	UNP B2HT54
C	0	SER	-	EXPRESSION TAG	UNP B2HT54
D	-3	GLY	-	EXPRESSION TAG	UNP B2HT54
D	-2	PRO	-	EXPRESSION TAG	UNP B2HT54
D	-1	GLY	-	EXPRESSION TAG	UNP B2HT54
D	0	SER	-	EXPRESSION TAG	UNP B2HT54

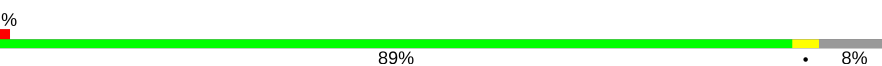
- Molecule 2 is water.

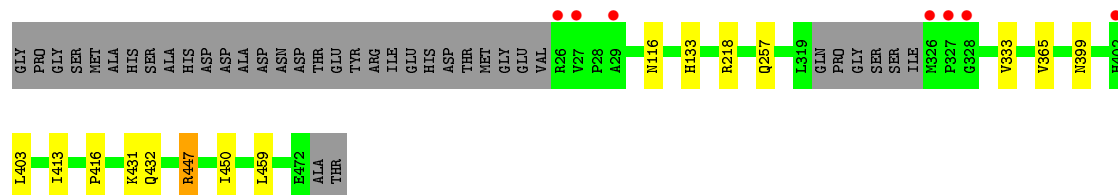
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	447	Total 447	O 447	0	0
2	B	435	Total 435	O 435	0	0
2	C	395	Total 395	O 395	0	0
2	D	407	Total 407	O 407	0	0

3 Residue-property plots [i](#)

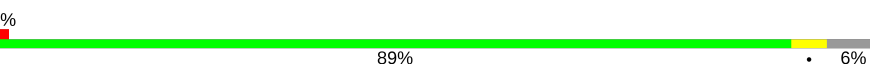
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are 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.

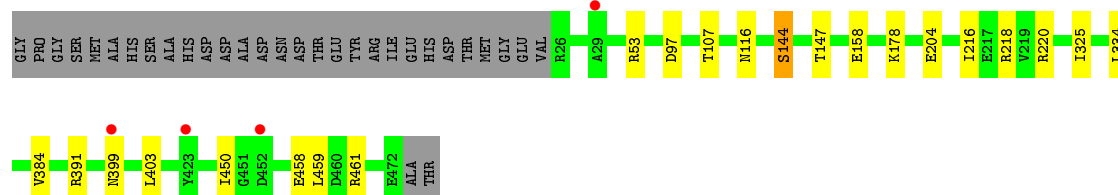
• Molecule 1: Fumarase Fum

Chain A: 




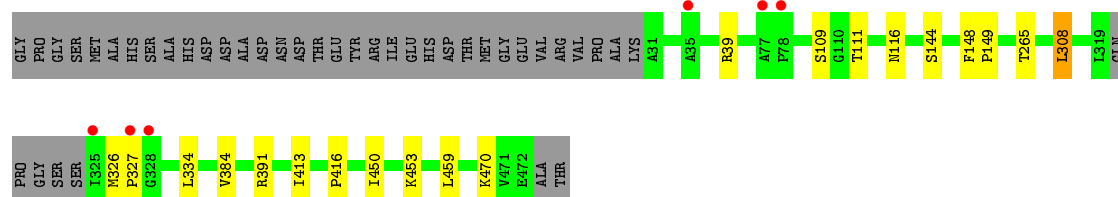
• Molecule 1: Fumarase Fum

Chain B: 




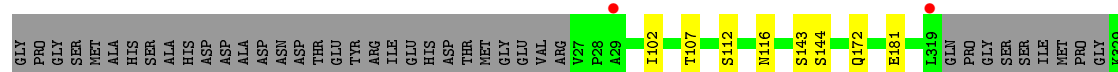
• Molecule 1: Fumarase Fum

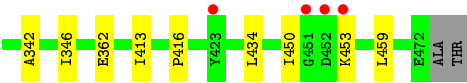
Chain C: 



• Molecule 1: Fumarase Fum

Chain D: 





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	97.88Å 114.08Å 98.84Å 90.00° 107.01° 90.00°	Depositor
Resolution (Å)	50.00 – 1.85 48.24 – 1.85	Depositor EDS
% Data completeness (in resolution range)	99.9 (50.00-1.85) 99.9 (48.24-1.85)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.18 (at 1.86Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.155 , 0.188 0.160 , 0.193	Depositor DCC
R_{free} test set	8834 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	15.1	Xtriage
Anisotropy	0.049	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 44.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.015 for l,-k,h	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	14756	wwPDB-VP
Average B, all atoms (Å ²)	17.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.27% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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	0.67	0/3347	0.63	1/4557 (0.0%)
1	B	0.70	0/3400	0.66	1/4627 (0.0%)
1	C	0.66	0/3373	0.64	0/4590
1	D	0.68	0/3267	0.62	0/4451
All	All	0.68	0/13387	0.64	2/18225 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	218	ARG	NE-CZ-NH2	-5.75	117.43	120.30
1	A	218	ARG	NE-CZ-NH2	-5.27	117.67	120.30

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	3269	0	3311	9	0
1	B	3315	0	3378	25	0
1	C	3282	0	3341	19	0
1	D	3206	0	3236	9	0
2	A	447	0	0	4	0
2	B	435	0	0	1	0
2	C	395	0	0	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	D	407	0	0	2	0
All	All	14756	0	13266	58	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 2.

All (58) 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:461[B]:ARG:NH2	1:B:461[B]:ARG:HB2	1.66	1.09
1:B:107:THR:HG21	2:C:879:HOH:O	1.52	1.08
1:C:111[B]:THR:HG22	1:C:144:SER:CB	1.87	1.05
1:C:39[B]:ARG:HE	1:C:39[B]:ARG:HA	1.20	1.00
1:C:326:MET:HG2	1:C:327:PRO:HD2	1.47	0.96
1:B:461[B]:ARG:HH21	1:B:461[B]:ARG:CG	1.86	0.88
1:B:461[B]:ARG:HH21	1:B:461[B]:ARG:HG3	1.42	0.85
1:B:461[B]:ARG:NH2	1:B:461[B]:ARG:CB	2.40	0.84
1:C:39[B]:ARG:CA	1:C:39[B]:ARG:HE	1.89	0.82
1:C:39[B]:ARG:NE	1:C:39[B]:ARG:HA	1.96	0.79
1:C:111[B]:THR:HG22	1:C:144:SER:OG	1.82	0.79
1:D:172:GLN:HG2	2:D:1061:HOH:O	1.83	0.78
1:C:265[B]:THR:HG21	2:C:1518:HOH:O	1.88	0.74
1:B:461[B]:ARG:CZ	1:B:461[B]:ARG:HB2	2.19	0.71
1:C:111[B]:THR:HG22	1:C:144:SER:HB2	1.70	0.70
1:C:109:SER:OG	1:C:111[B]:THR:HG23	1.93	0.69
1:B:461[B]:ARG:HH21	1:B:461[B]:ARG:CB	2.05	0.68
1:B:461[B]:ARG:NH2	1:B:461[B]:ARG:CG	2.53	0.66
1:C:111[B]:THR:HG21	2:C:1642:HOH:O	1.97	0.64
1:A:365:VAL:HG22	2:A:643:HOH:O	1.99	0.62
1:B:399:ASN:O	1:B:403:LEU:HG	2.00	0.61
1:A:257:GLN:OE1	2:A:1213:HOH:O	2.16	0.61
1:B:461[B]:ARG:CB	1:B:461[B]:ARG:CZ	2.78	0.59
1:C:111[B]:THR:CG2	1:C:144:SER:CB	2.75	0.57
2:A:687:HOH:O	1:C:470:LYS:HE3	2.06	0.55
1:D:102:ILE:HD13	1:D:112:SER:HB3	1.89	0.53
1:B:53[A]:ARG:NE	1:B:97:ASP:OD2	2.42	0.52
1:A:399:ASN:O	1:A:403:LEU:HG	2.10	0.52
1:D:413:ILE:O	1:D:416:PRO:HD2	2.13	0.48
1:A:432:GLN:OE1	1:A:447:ARG:NH1	2.44	0.48
1:C:334:LEU:HD11	1:C:391[B]:ARG:HB2	1.96	0.48
1:D:181[A]:GLU:HG3	2:D:1648:HOH:O	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:413:ILE:O	1:C:416:PRO:HD2	2.15	0.47
1:C:326:MET:CG	1:C:327:PRO:HD2	2.31	0.46
1:C:450:ILE:HD11	1:C:459:LEU:HD22	1.99	0.45
1:A:450:ILE:HD11	1:A:459:LEU:HD22	1.99	0.45
1:B:216:ILE:O	1:B:220[B]:ARG:HG3	2.17	0.45
1:B:178[B]:LYS:HD3	1:B:178[B]:LYS:HA	1.64	0.44
1:C:148:PHE:N	1:C:149:PRO:HD2	2.33	0.44
1:A:413:ILE:O	1:A:416:PRO:HD2	2.18	0.43
1:B:450:ILE:HD11	1:B:459:LEU:HD22	2.00	0.43
1:B:458:GLU:OE2	1:B:461[A]:ARG:NH2	2.41	0.43
1:B:334:LEU:HD11	1:B:391[B]:ARG:HB2	1.99	0.43
1:B:144:SER:HA	1:B:147:THR:HG1	1.84	0.42
1:D:342:ALA:O	1:D:346:ILE:HG12	2.19	0.42
1:A:133[A]:HIS:CE1	2:A:535:HOH:O	2.72	0.42
1:C:308[B]:LEU:HD21	1:D:434:LEU:HD22	2.01	0.42
1:A:333:VAL:CG1	1:D:107[A]:THR:HG22	2.49	0.41
1:B:204:GLU:OE2	1:D:362:GLU:OE2	2.38	0.41
1:C:334:LEU:CD1	1:C:391[B]:ARG:HB2	2.51	0.41
1:D:450:ILE:HD11	1:D:459:LEU:HD22	2.03	0.41
1:B:158:GLU:CG	2:B:678:HOH:O	2.69	0.40
1:A:431:LYS:HE3	1:B:325:ILE:O	2.22	0.40

There are no symmetry-related clashes.

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	448/478 (94%)	437 (98%)	11 (2%)	0	100	100
1	B	458/478 (96%)	451 (98%)	7 (2%)	0	100	100
1	C	450/478 (94%)	436 (97%)	14 (3%)	0	100	100
1	D	439/478 (92%)	429 (98%)	10 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	1795/1912 (94%)	1753 (98%)	42 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	338/366 (92%)	336 (99%)	2 (1%)	86	83
1	B	343/366 (94%)	340 (99%)	3 (1%)	78	72
1	C	341/366 (93%)	336 (98%)	5 (2%)	65	53
1	D	328/366 (90%)	324 (99%)	4 (1%)	71	62
All	All	1350/1464 (92%)	1336 (99%)	14 (1%)	76	69

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	116	ASN
1	A	447	ARG
1	B	116	ASN
1	B	144	SER
1	B	384	VAL
1	C	116	ASN
1	C	308[A]	LEU
1	C	308[B]	LEU
1	C	384	VAL
1	C	453	LYS
1	D	116	ASN
1	D	143	SER
1	D	144	SER
1	D	453	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	135	ASN
1	C	402	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, 95th 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 ²)	Q<0.9
1	A	441/478 (92%)	-0.53	7 (1%) 72 72	7, 14, 30, 45	0
1	B	447/478 (93%)	-0.51	4 (0%) 84 84	6, 14, 29, 46	0
1	C	437/478 (91%)	-0.38	6 (1%) 75 76	6, 16, 36, 51	0
1	D	437/478 (91%)	-0.48	6 (1%) 75 76	6, 15, 31, 45	0
All	All	1762/1912 (92%)	-0.47	23 (1%) 77 78	6, 14, 33, 51	0

All (23) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	452	ASP	5.4
1	A	328	GLY	5.2
1	C	325	ILE	4.5
1	D	452	ASP	4.3
1	B	29	ALA	3.3
1	D	451	GLY	3.0
1	D	29	ALA	3.0
1	C	35	ALA	2.7
1	C	77	ALA	2.6
1	A	26	ARG	2.5
1	D	319	LEU	2.4
1	A	326	MET	2.3
1	B	423	TYR	2.3
1	D	423	TYR	2.3
1	C	328	GLY	2.2
1	A	29	ALA	2.1
1	C	327	PRO	2.1
1	B	399	ASN	2.1
1	D	453	LYS	2.1
1	A	402	HIS	2.1
1	A	27	VAL	2.0

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Mol	Chain	Res	Type	RSRZ
1	A	327	PRO	2.0
1	C	78	PRO	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.