Full wwPDB X-ray Structure Validation Report (i)

Jun 18, 2024 – 02:08 AM EDT

PDB ID	:	5YM9
Title	:	Crystal Structure of the Deamidase from Legionella pneumophila
Authors	:	Zhu, Z.L.; Zhu, M.
Deposited on	:	2017-10-21
Resolution	:	2.50 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:

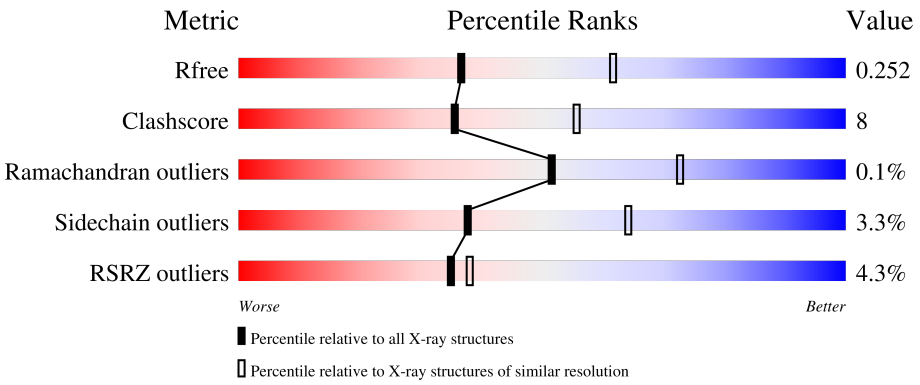
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.37.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

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 2.50 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	4661 (2.50-2.50)
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)
RSRZ outliers	127900	4559 (2.50-2.50)

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	384	4% 85% 13% •
1	B	384	4% 79% 19% •

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 6405 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 Uncharacterized protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	384	3095	1970	515	595	15	0	0	0
1	B	384	3095	1970	515	595	15	0	0	0

- Molecule 2 is water.

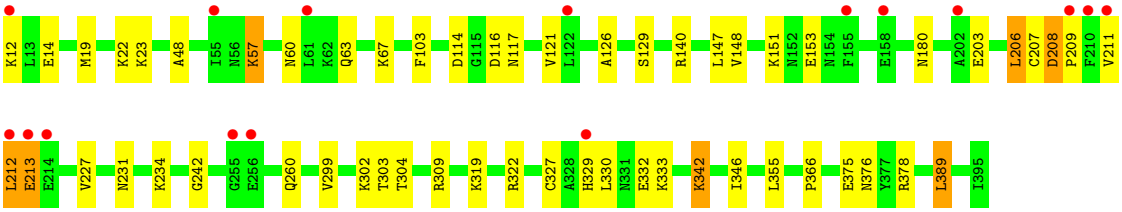
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
2	A	111	111	111	0	0
2	B	104	104	104	0	0

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.

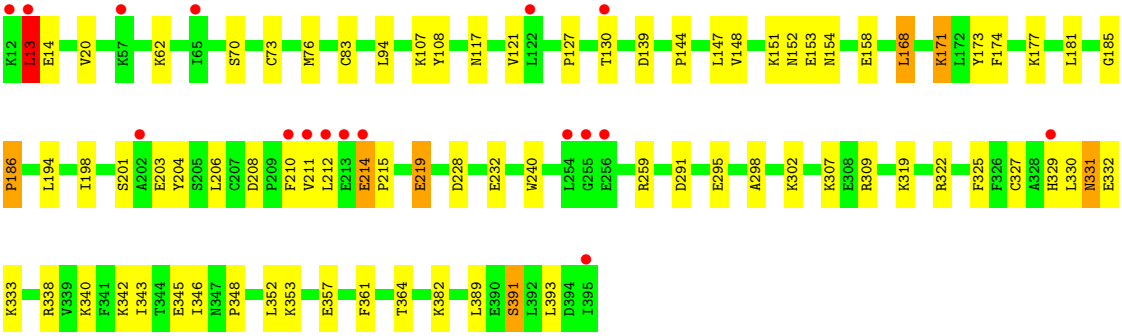
• Molecule 1: Uncharacterized protein

Chain A:



• Molecule 1: Uncharacterized protein

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α, β, γ	122.10Å 135.90Å 64.58Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.81 – 2.50 46.81 – 2.50	Depositor EDS
% Data completeness (in resolution range)	98.2 (46.81-2.50) 98.4 (46.81-2.50)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.78 (at 2.51Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.201 , 0.252 0.202 , 0.252	Depositor DCC
R_{free} test set	1804 reflections (4.83%)	wwPDB-VP
Wilson B-factor (Å^2)	59.4	Xtriage
Anisotropy	0.460	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.31 , 51.2	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	6405	wwPDB-VP
Average B, all atoms (Å^2)	70.0	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 37.80 % 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 4.0582e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

[1]Intensities estimated from amplitudes.

[2]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.

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	0.57	0/3161	0.69	2/4262 (0.0%)
1	B	0.62	2/3161 (0.1%)	0.73	5/4262 (0.1%)
All	All	0.59	2/6322 (0.0%)	0.71	7/8524 (0.1%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	186	PRO	N-CD	6.38	1.56	1.47
1	B	185	GLY	C-O	-5.83	1.14	1.23

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	210	PHE	N-CA-C	12.29	144.17	111.00
1	B	185	GLY	N-CA-C	8.60	134.61	113.10
1	B	210	PHE	C-N-CA	7.61	140.72	121.70
1	A	207	CYS	CB-CA-C	-5.86	98.68	110.40
1	B	186	PRO	CA-N-CD	-5.82	103.35	111.50
1	A	208	ASP	N-CA-CB	-5.78	100.20	110.60
1	B	13	LEU	CA-CB-CG	5.66	128.31	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	3095	0	3061	48	0
1	B	3095	0	3061	52	0
2	A	111	0	0	6	2
2	B	104	0	0	7	2
All	All	6405	0	6122	100	2

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 8.

All (100) 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:A:330:LEU:H	1:A:330:LEU:HD12	1.10	1.15
1:A:57:LYS:HD3	1:A:57:LYS:H	1.13	1.08
1:A:180:ASN:ND2	2:A:401:HOH:O	1.99	0.93
1:B:327:CYS:SG	1:B:333:LYS:NZ	2.44	0.89
1:A:330:LEU:H	1:A:330:LEU:CD1	1.85	0.88
1:A:330:LEU:HD12	1:A:330:LEU:N	1.90	0.86
1:B:83:CYS:SG	2:B:480:HOH:O	2.32	0.86
1:A:57:LYS:HD3	1:A:57:LYS:N	1.94	0.82
1:B:151:LYS:HB2	1:B:153:GLU:HG2	1.69	0.74
1:B:357:GLU:OE2	2:B:401:HOH:O	2.08	0.71
1:A:148:VAL:HG22	1:A:203:GLU:HG2	1.72	0.71
1:A:57:LYS:H	1:A:57:LYS:CD	1.94	0.69
1:B:382:LYS:NZ	2:B:405:HOH:O	2.25	0.69
1:A:140:ARG:NH1	2:A:403:HOH:O	2.26	0.68
1:B:307:LYS:NZ	2:B:406:HOH:O	2.26	0.68
1:A:147:LEU:HB2	1:A:206:LEU:HD21	1.76	0.68
1:A:304:THR:O	2:A:402:HOH:O	2.12	0.67
1:B:361:PHE:O	1:B:364:THR:OG1	2.13	0.67
1:B:291:ASP:O	1:B:295:GLU:HG3	1.95	0.66
1:B:20:VAL:HA	1:B:352:LEU:HD21	1.78	0.65
1:A:375:GLU:OE1	1:A:378:ARG:NH2	2.27	0.65
1:B:171:LYS:HG2	1:B:173:TYR:CZ	2.32	0.64
1:B:198:ILE:HA	1:B:219:GLU:HB2	1.79	0.63
1:B:208:ASP:OD2	1:B:211:VAL:HG21	1.98	0.63
1:A:12:LYS:NZ	1:A:22:LYS:NZ	2.47	0.62
1:A:12:LYS:NZ	1:A:22:LYS:HZ2	1.98	0.61
1:A:147:LEU:CB	1:A:206:LEU:HD21	2.31	0.60
1:A:366:PRO:HG2	1:A:376:ASN:ND2	2.17	0.60
1:B:177:LYS:NZ	1:B:240:TRP:O	2.25	0.60
1:A:327:CYS:SG	1:A:333:LYS:NZ	2.61	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:242:GLY:O	2:A:403:HOH:O	2.16	0.60
1:A:208:ASP:OD1	1:A:209:PRO:HD2	2.01	0.59
1:B:153:GLU:HG3	1:B:154:ASN:N	2.16	0.59
1:B:391:SER:OG	2:B:402:HOH:O	2.17	0.59
1:A:12:LYS:HZ2	1:A:22:LYS:NZ	2.01	0.59
1:B:228:ASP:O	1:B:232:GLU:HG2	2.04	0.57
1:B:214:GLU:N	1:B:215:PRO:HD3	2.19	0.57
1:B:329:HIS:O	1:B:332:GLU:HG3	2.05	0.56
1:A:19:MET:HE1	1:A:22:LYS:HD2	1.87	0.56
1:B:147:LEU:HB2	1:B:206:LEU:HD21	1.89	0.55
1:A:60:ASN:ND2	1:A:63:GLN:HE21	2.06	0.54
1:A:213:GLU:OE1	1:A:213:GLU:HA	2.06	0.54
1:B:117:ASN:O	1:B:121:VAL:HG23	2.07	0.54
1:A:208:ASP:O	1:A:211:VAL:HB	2.09	0.53
1:B:139:ASP:OD1	1:B:340:LYS:NZ	2.39	0.53
1:A:151:LYS:HG3	1:A:153:GLU:HG2	1.92	0.52
1:A:319:LYS:HD3	1:A:322:ARG:HE	1.76	0.50
1:B:298:ALA:O	1:B:302:LYS:HE2	2.11	0.50
1:A:60:ASN:HD22	1:A:63:GLN:HE21	1.60	0.50
1:B:201:SER:HB2	1:B:204:TYR:CZ	2.47	0.50
1:B:338:ARG:NH2	2:B:408:HOH:O	2.43	0.50
1:A:48:ALA:HB1	1:A:389:LEU:HD13	1.93	0.49
1:A:329:HIS:O	1:A:332:GLU:HG3	2.13	0.49
1:A:12:LYS:CE	1:A:22:LYS:HZ2	2.24	0.49
1:B:127:PRO:HA	1:B:259:ARG:HH21	1.78	0.49
1:B:144:PRO:HG3	1:B:174:PHE:CE1	2.47	0.49
1:A:329:HIS:N	1:A:329:HIS:CD2	2.81	0.49
1:A:103:PHE:CZ	1:A:342:LYS:HG2	2.48	0.48
1:A:212:LEU:HD21	1:A:227:VAL:HG11	1.95	0.48
1:A:14:GLU:OE1	2:A:404:HOH:O	2.18	0.48
1:A:23:LYS:HB3	1:A:355:LEU:HD13	1.96	0.48
1:B:319:LYS:HD3	1:B:322:ARG:HE	1.77	0.48
1:B:73:CYS:O	2:B:403:HOH:O	2.20	0.48
1:A:231:ASN:HA	1:A:234:LYS:HG3	1.95	0.47
1:B:151:LYS:CB	1:B:153:GLU:HG2	2.42	0.47
1:B:152:ASN:ND2	1:B:168:LEU:HD23	2.29	0.47
1:B:208:ASP:OD2	1:B:211:VAL:CG2	2.64	0.46
1:B:151:LYS:HB2	1:B:153:GLU:CG	2.44	0.46
1:B:214:GLU:N	1:B:215:PRO:CD	2.79	0.46
1:B:153:GLU:CG	1:B:154:ASN:N	2.78	0.46
1:B:346:ILE:O	1:B:348:PRO:HD3	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:148:VAL:HG22	1:B:203:GLU:HG2	1.98	0.46
1:B:107:LYS:HA	1:B:107:LYS:HD2	1.84	0.45
1:B:333:LYS:HD3	1:B:333:LYS:HA	1.73	0.45
1:A:126:ALA:HB3	1:A:129:SER:OG	2.17	0.45
1:B:322:ARG:HD3	1:B:325:PHE:HZ	1.81	0.45
1:A:114:ASP:HB2	1:A:116:ASP:OD1	2.17	0.45
1:B:389:LEU:HD23	1:B:389:LEU:HA	1.86	0.44
1:A:151:LYS:HB3	1:A:151:LYS:HE2	1.72	0.44
1:B:208:ASP:CG	1:B:211:VAL:HG23	2.38	0.44
1:A:208:ASP:OD1	1:A:209:PRO:CD	2.66	0.43
1:A:117:ASN:O	1:A:121:VAL:HG23	2.18	0.43
1:B:130:THR:CG2	1:B:345:GLU:HG2	2.48	0.43
1:B:194:LEU:O	1:B:198:ILE:HD12	2.19	0.43
1:B:331:ASN:HD22	1:B:331:ASN:C	2.21	0.43
1:A:309:ARG:HD2	1:A:333:LYS:O	2.19	0.43
1:B:208:ASP:CG	1:B:211:VAL:CG2	2.87	0.42
1:A:260:GLN:OE1	2:A:405:HOH:O	2.21	0.42
1:A:346:ILE:HD13	1:A:346:ILE:HA	1.88	0.42
1:A:23:LYS:HA	1:A:23:LYS:HD3	1.80	0.41
1:B:108:TYR:CZ	1:B:342:LYS:HE3	2.55	0.41
1:A:299:VAL:O	1:A:303:THR:HG22	2.20	0.41
1:A:67:LYS:HB3	1:A:67:LYS:HE2	1.88	0.41
1:B:20:VAL:CA	1:B:352:LEU:HD21	2.50	0.41
1:B:173:TYR:CG	1:B:181:LEU:HD11	2.55	0.41
1:B:13:LEU:HD22	1:B:14:GLU:N	2.36	0.41
1:A:333:LYS:HA	1:A:333:LYS:HD3	1.90	0.40
1:B:309:ARG:HD2	1:B:333:LYS:O	2.20	0.40
1:B:346:ILE:HD13	1:B:346:ILE:HA	1.84	0.40
1:B:393:LEU:HD23	1:B:393:LEU:HA	1.87	0.40

All (2) 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)
2:A:463:HOH:O	2:B:490:HOH:O[2_565]	2.13	0.07
2:A:509:HOH:O	2:B:502:HOH:O[4_455]	2.15	0.05

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	382/384 (100%)	365 (96%)	17 (4%)	0	100	100
1	B	382/384 (100%)	364 (95%)	17 (4%)	1 (0%)	41	61
All	All	764/768 (100%)	729 (95%)	34 (4%)	1 (0%)	51	73

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	214	GLU

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	345/345 (100%)	338 (98%)	7 (2%)	55	79
1	B	345/345 (100%)	329 (95%)	16 (5%)	27	50
All	All	690/690 (100%)	667 (97%)	23 (3%)	38	64

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	57	LYS
1	A	206	LEU
1	A	212	LEU
1	A	213	GLU
1	A	302	LYS

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Mol	Chain	Res	Type
1	A	342	LYS
1	A	389	LEU
1	B	13	LEU
1	B	62	LYS
1	B	70	SER
1	B	76	MET
1	B	94	LEU
1	B	158	GLU
1	B	168	LEU
1	B	171	LYS
1	B	186	PRO
1	B	212	LEU
1	B	219	GLU
1	B	330	LEU
1	B	331	ASN
1	B	343	ILE
1	B	353	LYS
1	B	391	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	47	GLN
1	A	63	GLN
1	A	180	ASN
1	A	329	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 monosaccharides 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 (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, 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	384/384 (100%)	0.34	16 (4%)	36	39	44, 68, 98, 133	0
1	B	384/384 (100%)	0.35	17 (4%)	34	37	45, 70, 100, 128	0
All	All	768/768 (100%)	0.35	33 (4%)	35	38	44, 69, 100, 133	0

All (33) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	210	PHE	8.2
1	A	212	LEU	6.6
1	B	212	LEU	6.4
1	B	210	PHE	5.9
1	A	209	PRO	5.3
1	B	12	LYS	5.0
1	B	211	VAL	4.8
1	B	256	GLU	4.1
1	B	213	GLU	3.5
1	A	214	GLU	3.3
1	A	211	VAL	3.2
1	A	155	PHE	3.1
1	B	202	ALA	3.1
1	A	12	LYS	3.1
1	A	122	LEU	2.9
1	B	329	HIS	2.9
1	A	158	GLU	2.8
1	A	255	GLY	2.8
1	B	13	LEU	2.8
1	B	214	GLU	2.7
1	B	57	LYS	2.6
1	B	130	THR	2.6
1	A	213	GLU	2.5
1	A	61	LEU	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	122	LEU	2.5
1	B	254	LEU	2.4
1	A	256	GLU	2.4
1	A	329	HIS	2.4
1	B	65	ILE	2.3
1	A	202	ALA	2.2
1	B	255	GLY	2.1
1	A	55	ILE	2.1
1	B	395	ILE	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)

There are no ligands in this entry.

6.5 Other polymers (i)

There are no such residues in this entry.