



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:04 AM GMT

PDB ID : 2ZD7  
Title : The structure of VPS75 (Vacuolar protein sorting-associated protein 75)  
Authors : Park, Y.J.; Luger, K.  
Deposited on : 2007-11-20  
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](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

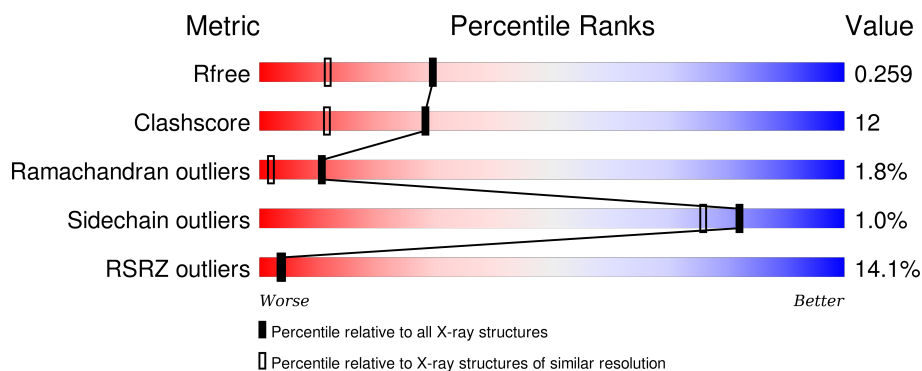
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 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))
$R_{free}$	91344	1745 (1.86-1.86)
Clashscore	102246	1898 (1.86-1.86)
Ramachandran outliers	100387	1875 (1.86-1.86)
Sidechain outliers	100360	1875 (1.86-1.86)
RSRZ outliers	91569	1747 (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  $\geq 3$ , 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	264	<div> <div>10%</div> <div> <div></div> <div>65%</div> <div>17%</div> <div>••</div> <div>16%</div> </div> </div>
1	B	264	<div> <div>11%</div> <div> <div></div> <div>63%</div> <div>16%</div> <div>•</div> <div>20%</div> </div> </div>
2	C	13	<div> <div>54%</div> <div> <div></div> <div>69%</div> <div>23%</div> <div>8%</div> </div> </div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3991 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 Vacuolar protein sorting-associated protein 75.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	222	Total	C	N	O	S	0	0	0
			1853	1191	301	356	5			
1	B	212	Total	C	N	O	S	0	0	0
			1773	1148	286	334	5			

- Molecule 2 is a protein called EVDLPLSDEEPSS.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	13	Total	C	N	O	0	0	0
			99	59	13	27			

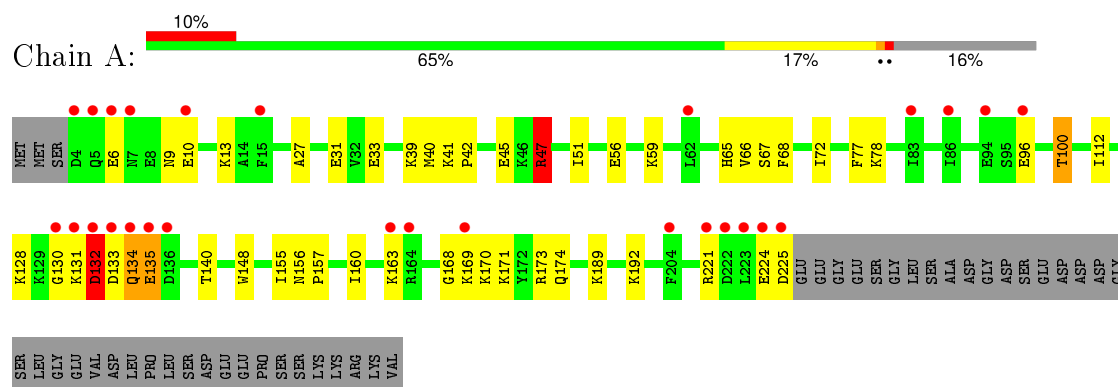
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	156	Total	O	0	0
			156	156		
3	B	110	Total	O	0	0
			110	110		

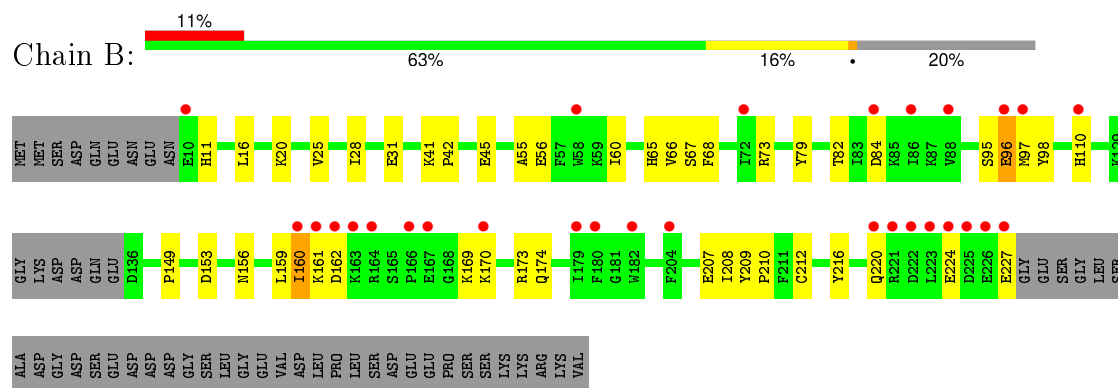
### 3 Residue-property plots [i](#)

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

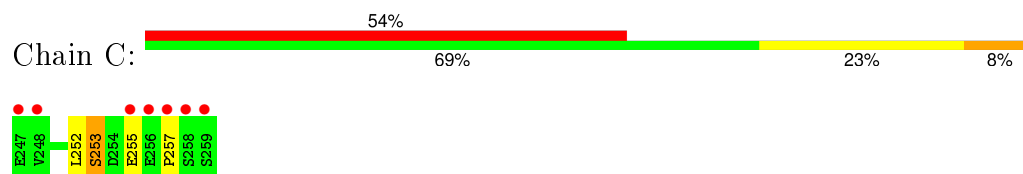
- Molecule 1: Vacuolar protein sorting-associated protein 75



- Molecule 1: Vacuolar protein sorting-associated protein 75



- Molecule 2: EVDLPLSDEEPSS



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	80.69Å 84.58Å 86.22Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 1.85 27.65 – 1.85	Depositor EDS
% Data completeness (in resolution range)	95.3 (50.00-1.85) 94.7 (27.65-1.85)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.31 (at 1.85Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.232 , 0.255 0.237 , 0.259	Depositor DCC
$R_{free}$ test set	4886 reflections (10.09%)	DCC
Wilson B-factor (Å <sup>2</sup> )	32.8	Xtriage
Anisotropy	0.334	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 43.8	EDS
Estimated twinning fraction	0.005 for -h,l,k	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 49841 reflections	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	3991	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.33% 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	0.36	0/1901	0.57	1/2561 (0.0%)
1	B	0.34	0/1820	0.55	0/2451
2	C	0.32	0/100	0.49	0/135
All	All	0.35	0/3821	0.56	1/5147 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	47	ARG	NE-CZ-NH2	-5.64	117.48	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	1853	0	1781	46	0
1	B	1773	0	1718	41	0
2	C	99	0	85	2	0
3	A	156	0	0	4	0
3	B	110	0	0	1	0
All	All	3991	0	3584	85	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 12.

All (85) 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:66:VAL:HG13	1:B:67:SER:H	1.40	0.86
1:B:79:TYR:O	1:B:82:THR:HG22	1.77	0.85
1:B:65:HIS:HD2	1:B:68:PHE:H	1.25	0.83
1:B:170:LYS:HG2	1:B:174:GLN:HE21	1.48	0.79
1:A:33:GLU:HG2	1:A:100:THR:CG2	2.14	0.77
1:A:33:GLU:HG2	1:A:100:THR:HG21	1.65	0.76
1:B:96:GLU:C	1:B:98:TYR:H	1.93	0.72
1:A:27:ALA:O	1:A:31:GLU:HG3	1.91	0.71
1:A:65:HIS:HD2	1:A:68:PHE:H	1.39	0.71
1:B:66:VAL:HG13	1:B:67:SER:N	2.06	0.70
1:B:84:ASP:OD2	1:B:110:HIS:HB2	1.92	0.70
1:B:41:LYS:HB3	1:B:42:PRO:HD3	1.76	0.67
1:A:41:LYS:HB3	1:A:42:PRO:HD3	1.76	0.66
1:A:40:MET:HE1	1:B:25:VAL:HG22	1.77	0.64
1:A:39:LYS:HE3	1:B:31:GLU:OE1	1.97	0.64
1:A:157:PRO:HG3	3:A:312:HOH:O	1.98	0.63
1:A:224:GLU:O	1:A:225:ASP:HB3	1.97	0.63
1:A:133:ASP:HB3	1:A:135:GLU:HG3	1.81	0.62
1:A:40:MET:CE	1:B:25:VAL:HG22	2.29	0.62
1:A:72:ILE:HD11	1:A:77:PHE:CE1	2.35	0.62
1:A:39:LYS:HE2	3:A:335:HOH:O	1.99	0.61
1:B:96:GLU:O	1:B:98:TYR:N	2.33	0.61
1:A:39:LYS:HD3	1:B:28:ILE:HG12	1.82	0.61
1:B:156:ASN:O	1:B:160:ILE:HG13	2.01	0.60
1:B:41:LYS:O	1:B:45:GLU:HG3	2.02	0.60
1:A:156:ASN:O	1:A:160:ILE:HG13	2.02	0.60
1:A:66:VAL:HG22	3:A:304:HOH:O	2.04	0.57
1:B:169:LYS:O	1:B:173:ARG:HG3	2.05	0.56
1:A:47:ARG:HD3	1:A:51:ILE:HD12	1.88	0.56
1:A:41:LYS:O	1:A:45:GLU:HG3	2.05	0.56
1:B:56:GLU:O	1:B:60:ILE:HG13	2.06	0.55
1:B:11:HIS:HA	3:B:322:HOH:O	2.07	0.55
1:A:221:ARG:O	1:A:224:GLU:HG2	2.07	0.54
1:A:6:GLU:O	1:A:10:GLU:HG2	2.08	0.54
1:A:192:LYS:O	1:A:192:LYS:HG2	2.09	0.53
1:A:33:GLU:HG2	1:A:100:THR:HG23	1.90	0.53
1:A:163:LYS:HA	1:A:168:GLY:C	2.30	0.53
1:A:47:ARG:HD3	1:A:51:ILE:CD1	2.39	0.52
1:B:73:ARG:HG2	1:B:73:ARG:HH11	1.73	0.52
1:A:100:THR:HG22	3:A:393:HOH:O	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:130:GLY:O	1:A:131:LYS:HB2	2.10	0.51
1:B:60:ILE:HD12	1:B:216:TYR:HE1	1.76	0.51
1:B:224:GLU:HA	1:B:227:GLU:HG3	1.93	0.50
2:C:255:GLU:C	2:C:257:PRO:HD3	2.32	0.50
1:A:131:LYS:O	1:A:132:ASP:HB2	2.12	0.49
1:A:9:ASN:ND2	1:A:13:LYS:HE3	2.28	0.49
1:B:224:GLU:HA	1:B:227:GLU:CD	2.33	0.49
1:B:96:GLU:C	1:B:98:TYR:N	2.60	0.48
1:A:78:LYS:HE2	1:A:112:ILE:HD13	1.95	0.48
1:B:73:ARG:NH1	1:B:73:ARG:HG2	2.29	0.48
1:B:16:LEU:HD23	1:B:16:LEU:C	2.34	0.48
1:B:95:SER:C	1:B:96:GLU:O	2.50	0.48
1:B:65:HIS:HE1	1:B:207:GLU:OE1	1.98	0.47
1:B:66:VAL:CG1	1:B:67:SER:H	2.18	0.47
1:A:128:LYS:HD2	1:A:140:THR:HG21	1.97	0.47
1:B:209:TYR:HB3	1:B:210:PRO:HD3	1.95	0.47
1:A:160:ILE:HG21	1:A:163:LYS:HB3	1.97	0.47
1:A:171:LYS:HE2	1:A:171:LYS:HA	1.97	0.46
1:A:148:TRP:CH2	1:A:157:PRO:HD3	2.51	0.46
1:B:220:GLN:O	1:B:224:GLU:HB2	2.16	0.46
1:B:65:HIS:HD2	1:B:68:PHE:N	2.02	0.46
1:A:189:LYS:N	1:A:189:LYS:HD2	2.32	0.45
1:B:79:TYR:CE2	1:B:149:PRO:HB3	2.51	0.45
2:C:252:LEU:O	2:C:253:SER:HB3	2.16	0.45
1:B:153:ASP:HA	1:B:159:LEU:HD12	1.98	0.45
1:B:224:GLU:HA	1:B:227:GLU:CG	2.47	0.44
1:A:170:LYS:HG2	1:A:174:GLN:HE21	1.82	0.44
1:A:224:GLU:O	1:A:225:ASP:CB	2.65	0.44
1:A:40:MET:HA	1:A:40:MET:HE3	1.99	0.43
1:A:169:LYS:O	1:A:173:ARG:HG3	2.18	0.43
1:A:65:HIS:CD2	1:A:67:SER:H	2.37	0.43
1:A:56:GLU:OE1	1:A:59:LYS:CE	2.66	0.43
1:B:170:LYS:HG2	1:B:174:GLN:NE2	2.26	0.42
1:A:134:GLN:O	1:A:135:GLU:C	2.57	0.42
1:A:9:ASN:O	1:A:13:LYS:HG3	2.19	0.42
1:B:161:LYS:O	1:B:162:ASP:HB2	2.19	0.42
1:A:160:ILE:CG2	1:A:163:LYS:HB3	2.48	0.42
1:B:159:LEU:O	1:B:160:ILE:C	2.59	0.42
1:A:155:ILE:HG22	1:A:171:LYS:HE2	2.02	0.42
1:A:65:HIS:CD2	1:A:68:PHE:H	2.28	0.41
1:B:95:SER:O	1:B:96:GLU:O	2.38	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:208:ILE:O	1:B:212:CYS:HB3	2.21	0.41
1:B:170:LYS:O	1:B:174:GLN:HG3	2.21	0.40
1:A:72:ILE:O	1:A:72:ILE:HD12	2.21	0.40
1:B:16:LEU:HD21	1:B:20:LYS:HE3	2.03	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	220/264 (83%)	208 (94%)	9 (4%)	3 (1%)	14	3
1	B	208/264 (79%)	196 (94%)	8 (4%)	4 (2%)	10	2
2	C	11/13 (85%)	10 (91%)	0	1 (9%)	1	0
All	All	439/541 (81%)	414 (94%)	17 (4%)	8 (2%)	11	2

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	135	GLU
1	B	55	ALA
1	B	96	GLU
1	A	132	ASP
1	B	160	ILE
2	C	253	SER
1	A	134	GLN
1	B	97	MET

### 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	199/235 (85%)	195 (98%)	4 (2%)	63	47
1	B	190/235 (81%)	190 (100%)	0	100	100
2	C	13/13 (100%)	13 (100%)	0	100	100
All	All	402/483 (83%)	398 (99%)	4 (1%)	82	76

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	47	ARG
1	A	96	GLU
1	A	100	THR
1	A	132	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	9	ASN
1	A	65	HIS
1	A	174	GLN
1	B	65	HIS
1	B	174	GLN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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	222/264 (84%)	0.77	27 (12%) 5 5	23, 34, 77, 84	0
1	B	212/264 (80%)	0.73	29 (13%) 4 4	26, 39, 68, 91	0
2	C	13/13 (100%)	4.04	7 (53%) 0 0	47, 67, 81, 82	0
All	All	447/541 (82%)	0.84	63 (14%) 4 3	23, 37, 77, 91	0

All (63) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	C	259	SER	12.8
1	A	133	ASP	12.4
1	A	132	ASP	12.1
1	A	134	GLN	9.4
2	C	257	PRO	7.6
2	C	258	SER	7.5
1	B	223	LEU	6.7
1	B	164	ARG	6.4
2	C	256	GLU	6.3
1	A	163	LYS	5.8
1	B	225	ASP	5.8
1	B	227	GLU	5.6
1	A	131	LYS	5.5
1	A	164	ARG	5.4
1	B	162	ASP	5.3
1	A	4	ASP	5.0
1	B	224	GLU	4.9
1	A	5	GLN	4.8
1	A	7	ASN	4.8
1	B	160	ILE	4.8
2	C	247	GLU	4.8
2	C	255	GLU	4.6
1	B	166	PRO	4.5

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Mol	Chain	Res	Type	RSRZ
1	B	226	GLU	4.4
1	A	225	ASP	4.3
1	B	221	ARG	4.1
1	A	221	ARG	4.0
1	B	167	GLU	3.6
1	B	179	ILE	3.6
1	B	161	LYS	3.5
1	B	222	ASP	3.5
2	C	248	VAL	3.3
1	A	223	LEU	3.2
1	B	204	PHE	3.1
1	A	136	ASP	3.1
1	A	6	GLU	3.0
1	B	163	LYS	2.9
1	A	96	GLU	2.9
1	B	10	GLU	2.8
1	B	96	GLU	2.8
1	B	58	TRP	2.8
1	B	88	VAL	2.6
1	A	204	PHE	2.6
1	B	97	MET	2.5
1	A	130	GLY	2.3
1	A	15	PHE	2.3
1	A	224	GLU	2.3
1	B	110	HIS	2.3
1	A	86	ILE	2.3
1	A	62	LEU	2.3
1	B	86	ILE	2.3
1	B	84	ASP	2.3
1	A	135	GLU	2.3
1	B	72	ILE	2.2
1	B	220	GLN	2.2
1	A	169	LYS	2.1
1	A	222	ASP	2.1
1	A	94	GLU	2.1
1	B	182	TRP	2.1
1	B	170	LYS	2.1
1	A	10	GLU	2.0
1	B	180	PHE	2.0
1	A	83	ILE	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.