



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 11:54 AM GMT

PDB ID : 3QAC  
Title : Structure of amaranth 11S proglobulin seed storage protein from *Amaranthus hypochondriacus* L.  
Authors : Tandang-Silvas, M.R.; Carrazco-Pena, L.; Barba de la Rosa, A.P.; Osuna-Castro, J.A.; Utsumi, S.; Mikami, B.; Maruyama, N.  
Deposited on : 2011-01-10  
Resolution : 2.27 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

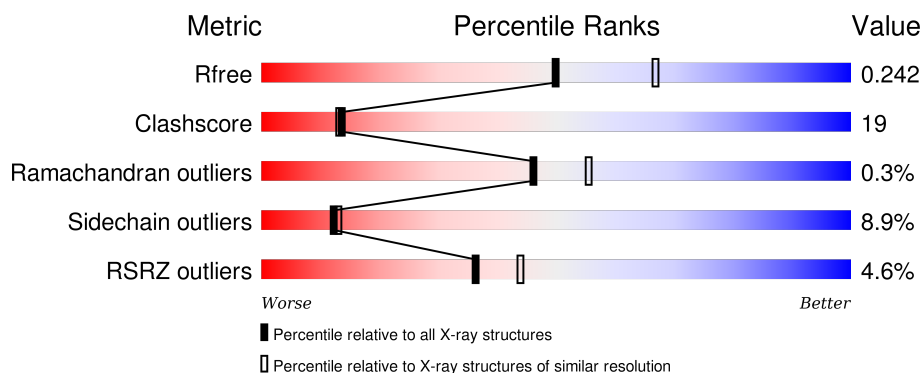
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*


The reported resolution of this entry is 2.27 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	5193 (2.30-2.26)
Clashscore	102246	5929 (2.30-2.26)
Ramachandran outliers	100387	5851 (2.30-2.26)
Sidechain outliers	100360	5850 (2.30-2.26)
RSRZ outliers	91569	5204 (2.30-2.26)

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	465	

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3027 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 11S globulin seed storage protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	372	Total	C	N	O	S	11	0	0
			2927	1830	529	556	12			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	EXPRESSION TAG	UNP Q38712

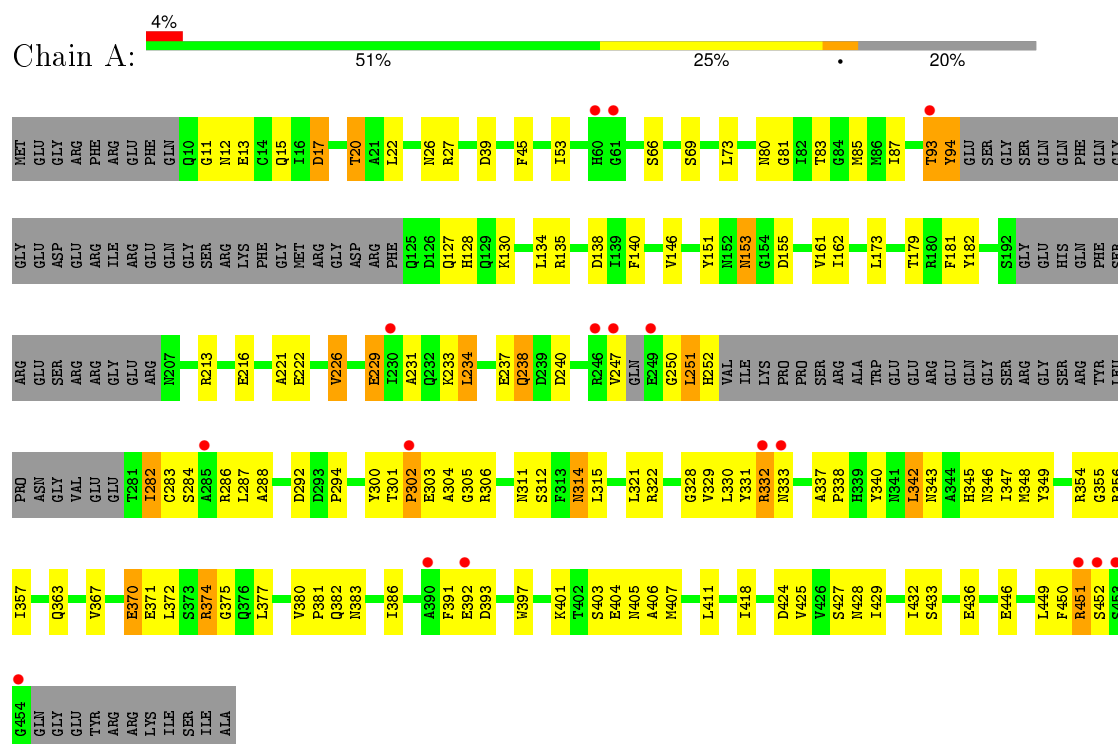
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	100	Total	O	0	0
			100	100		

### 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: 11S globulin seed storage protein



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 63	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	94.08 Å 94.08 Å 74.67 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	28.47 – 2.27 39.80 – 2.28	Depositor EDS
% Data completeness (in resolution range)	98.1 (28.47-2.27) 98.1 (39.80-2.28)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.43 (at 2.27 Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.5_2)	Depositor
R, $R_{free}$	0.188 , 0.249 0.183 , 0.242	Depositor DCC
$R_{free}$ test set	863 reflections (5.07%)	DCC
Wilson B-factor (Å <sup>2</sup> )	43.9	Xtriage
Anisotropy	0.362	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 45.2	EDS
Estimated twinning fraction	0.045 for h,-h-k,-l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 17038 reflections	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	3027	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.97% 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	0.46	0/2982	0.60	0/4035

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	2927	0	2847	107	0
2	A	100	0	0	2	0
All	All	3027	0	2847	107	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 19.

All (107) 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:374:ARG:HD2	1:A:375:GLY:N	1.72	1.03
1:A:311:ASN:HB2	1:A:405:ASN:HD22	1.32	0.90
1:A:300:TYR:OH	1:A:302:PRO:HA	1.73	0.88
1:A:374:ARG:C	1:A:374:ARG:HD2	2.06	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:251:LEU:O	1:A:252:HIS:HB2	1.91	0.69
1:A:85:MET:HE1	1:A:140:PHE:CE2	2.29	0.67
1:A:251:LEU:HD12	1:A:251:LEU:N	2.10	0.66
1:A:300:TYR:CZ	1:A:302:PRO:HA	2.33	0.64
1:A:17:ASP:HA	1:A:367:VAL:HG12	1.80	0.63
1:A:11:GLY:HA2	1:A:13:GLU:OE2	2.00	0.61
1:A:370:GLU:HG2	2:A:488:HOH:O	2.02	0.59
1:A:20:THR:HB	2:A:505:HOH:O	2.02	0.58
1:A:151:TYR:CE2	1:A:251:LEU:HD13	2.38	0.58
1:A:337:ALA:HB2	1:A:450:PHE:CE1	2.39	0.57
1:A:314:ASN:HD22	1:A:315:LEU:N	2.02	0.57
1:A:312:SER:H	1:A:405:ASN:ND2	2.03	0.57
1:A:321:LEU:N	1:A:321:LEU:HD23	2.20	0.57
1:A:333:ASN:O	1:A:452:SER:HB2	2.04	0.56
1:A:13:GLU:CD	1:A:13:GLU:H	2.09	0.55
1:A:238:GLN:HE21	1:A:238:GLN:HA	1.71	0.55
1:A:130:LYS:NZ	1:A:284:SER:HA	2.21	0.55
1:A:332:ARG:HD2	1:A:451:ARG:NH2	2.23	0.54
1:A:93:THR:HB	1:A:128:HIS:O	2.07	0.54
1:A:451:ARG:HD2	1:A:452:SER:H	1.72	0.54
1:A:311:ASN:HB2	1:A:405:ASN:ND2	2.13	0.54
1:A:425:VAL:O	1:A:429:ILE:HG12	2.08	0.54
1:A:53:ILE:HG13	1:A:162:ILE:HB	1.90	0.54
1:A:451:ARG:HD2	1:A:452:SER:N	2.24	0.53
1:A:282:ILE:HG12	1:A:283:CYS:N	2.23	0.53
1:A:355:GLY:O	1:A:371:GLU:HA	2.09	0.53
1:A:338:PRO:HB2	1:A:411:LEU:HD12	1.91	0.53
1:A:161:VAL:HG21	1:A:349:TYR:CE2	2.44	0.52
1:A:330:LEU:HD21	1:A:449:LEU:HD22	1.90	0.52
1:A:347:ILE:HD12	1:A:347:ILE:N	2.25	0.52
1:A:377:LEU:C	1:A:377:LEU:HD12	2.31	0.51
1:A:301:THR:HB	1:A:304:ALA:HB3	1.91	0.51
1:A:135:ARG:O	1:A:138:ASP:HB2	2.10	0.51
1:A:81:GLY:HA2	1:A:153:ASN:ND2	2.25	0.51
1:A:348:MET:CE	1:A:357:ILE:HD13	2.42	0.50
1:A:404:GLU:O	1:A:405:ASN:C	2.50	0.49
1:A:305:GLY:HA2	1:A:331:TYR:CE2	2.48	0.49
1:A:401:LYS:HD3	1:A:406:ALA:HB3	1.94	0.49
1:A:286:ARG:O	1:A:287:LEU:HB2	2.13	0.48
1:A:346:ASN:HB3	1:A:380:VAL:HB	1.96	0.48
1:A:179:THR:HG22	1:A:181:PHE:CE2	2.49	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:17:ASP:HA	1:A:367:VAL:CG1	2.44	0.48
1:A:26:ASN:OD1	1:A:27:ARG:N	2.47	0.48
1:A:83:THR:HB	1:A:134:LEU:HD11	1.95	0.48
1:A:130:LYS:HZ3	1:A:284:SER:HA	1.79	0.47
1:A:392:GLU:H	1:A:392:GLU:CD	2.18	0.47
1:A:357:ILE:HD12	1:A:372:LEU:HB2	1.96	0.47
1:A:85:MET:CE	1:A:140:PHE:CE2	2.97	0.47
1:A:69:SER:HB3	1:A:173:LEU:HD12	1.97	0.47
1:A:151:TYR:CZ	1:A:251:LEU:CD1	2.98	0.46
1:A:328:GLY:HA3	1:A:397:TRP:CZ2	2.51	0.46
1:A:311:ASN:H	1:A:314:ASN:HD21	1.63	0.46
1:A:94:TYR:CD1	1:A:127:GLN:HG2	2.51	0.46
1:A:312:SER:H	1:A:405:ASN:HD21	1.63	0.46
1:A:312:SER:HB2	1:A:405:ASN:HD21	1.80	0.46
1:A:85:MET:HG2	1:A:146:VAL:HG11	1.97	0.46
1:A:231:ALA:O	1:A:234:LEU:HB2	2.16	0.45
1:A:140:PHE:CD1	1:A:288:ALA:HB2	2.52	0.45
1:A:348:MET:HE1	1:A:357:ILE:HD13	1.97	0.45
1:A:427:SER:HB2	1:A:432:ILE:O	2.16	0.45
1:A:87:ILE:HD13	1:A:282:ILE:HG13	1.98	0.45
1:A:93:THR:O	1:A:127:GLN:HA	2.17	0.45
1:A:20:THR:HG22	1:A:22:LEU:HD23	1.99	0.45
1:A:73:LEU:HA	1:A:73:LEU:HD12	1.84	0.45
1:A:12:ASN:HD22	1:A:15:GLN:HG3	1.81	0.45
1:A:433:SER:OG	1:A:436:GLU:HG3	2.17	0.45
1:A:301:THR:HG21	1:A:446:GLU:HG2	1.98	0.45
1:A:424:ASP:O	1:A:428:ASN:HB2	2.16	0.45
1:A:371:GLU:HG2	1:A:391:PHE:HE2	1.82	0.45
1:A:305:GLY:HA3	1:A:329:VAL:O	2.17	0.45
1:A:151:TYR:CZ	1:A:251:LEU:HD13	2.52	0.44
1:A:300:TYR:HH	1:A:302:PRO:HA	1.77	0.44
1:A:213:ARG:HG3	1:A:213:ARG:O	2.17	0.44
1:A:303:GLU:O	1:A:451:ARG:HD3	2.18	0.44
1:A:357:ILE:HD12	1:A:372:LEU:HG	2.00	0.44
1:A:234:LEU:HA	1:A:234:LEU:HD12	1.85	0.44
1:A:382:GLN:O	1:A:383:ASN:HB2	2.16	0.44
1:A:251:LEU:O	1:A:252:HIS:CB	2.64	0.43
1:A:85:MET:CE	1:A:140:PHE:CZ	3.01	0.43
1:A:411:LEU:O	1:A:418:ILE:HG12	2.17	0.43
1:A:151:TYR:CE1	1:A:251:LEU:HD11	2.53	0.43
1:A:337:ALA:HB2	1:A:450:PHE:CZ	2.54	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:69:SER:CB	1:A:173:LEU:HD12	2.49	0.43
1:A:340:TYR:CE2	1:A:342:LEU:HD22	2.54	0.43
1:A:240:ASP:OD2	1:A:240:ASP:N	2.50	0.43
1:A:332:ARG:HD2	1:A:451:ARG:HH22	1.84	0.43
1:A:153:ASN:C	1:A:153:ASN:HD22	2.22	0.42
1:A:322:ARG:O	1:A:403:SER:HA	2.19	0.42
1:A:404:GLU:HG2	1:A:405:ASN:OD1	2.19	0.42
1:A:221:ALA:HB1	1:A:226:VAL:O	2.20	0.42
1:A:314:ASN:C	1:A:314:ASN:HD22	2.23	0.42
1:A:354:ARG:HD3	1:A:371:GLU:OE1	2.20	0.42
1:A:292:ASP:O	1:A:294:PRO:HD3	2.19	0.42
1:A:250:GLY:C	1:A:251:LEU:HD12	2.39	0.42
1:A:338:PRO:HA	1:A:386:ILE:O	2.20	0.41
1:A:233:LYS:HA	1:A:233:LYS:HD2	1.77	0.41
1:A:322:ARG:HA	1:A:404:GLU:HB2	2.03	0.41
1:A:39:ASP:O	1:A:45:PHE:HD1	2.04	0.41
1:A:306:ARG:HG2	1:A:329:VAL:HG22	2.03	0.41
1:A:229:GLU:H	1:A:229:GLU:HG2	1.66	0.41
1:A:345:HIS:ND1	1:A:381:PRO:HA	2.35	0.41
1:A:238:GLN:HA	1:A:238:GLN:NE2	2.35	0.40
1:A:66:SER:HB3	1:A:182:TYR:CD2	2.56	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	362/465 (78%)	343 (95%)	18 (5%)	1 (0%)	46 55

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	302	PRO

### 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	315/394 (80%)	287 (91%)	28 (9%)	12	13

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	17	ASP
1	A	20	THR
1	A	80	ASN
1	A	93	THR
1	A	94	TYR
1	A	153	ASN
1	A	155	ASP
1	A	216	GLU
1	A	222	GLU
1	A	226	VAL
1	A	229	GLU
1	A	234	LEU
1	A	237	GLU
1	A	238	GLN
1	A	247	VAL
1	A	251	LEU
1	A	282	ILE
1	A	314	ASN
1	A	332	ARG
1	A	342	LEU
1	A	343	ASN
1	A	356	ARG
1	A	363	GLN
1	A	370	GLU
1	A	374	ARG
1	A	393	ASP

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Mol	Chain	Res	Type
1	A	407	MET
1	A	451	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	12	ASN
1	A	125	GLN
1	A	153	ASN
1	A	172	GLN
1	A	232	GLN
1	A	238	GLN
1	A	314	ASN
1	A	382	GLN
1	A	383	ASN
1	A	405	ASN
1	A	409	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

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, 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	372/465 (80%)	0.13	17 (4%) 36 44	31, 48, 71, 84	5 (1%)

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	453	SER	7.1
1	A	454	GLY	6.4
1	A	452	SER	4.2
1	A	246	ARG	3.1
1	A	249	GLU	3.0
1	A	61	GLY	2.9
1	A	332	ARG	2.9
1	A	247	VAL	2.6
1	A	392	GLU	2.6
1	A	230	ILE	2.5
1	A	60	HIS	2.4
1	A	451	ARG	2.3
1	A	302	PRO	2.3
1	A	285	ALA	2.1
1	A	390	ALA	2.1
1	A	93	THR	2.1
1	A	333	ASN	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 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.