



# wwPDB X-ray Structure Validation Summary Report ⓘ

May 13, 2020 – 06:43 am BST

PDB ID : 2JCC  
Title : AH3 recognition of mutant HLA-A2 W167A  
Authors : Miller, P.; Benhar, Y.P.; Biddison, W.; Collins, E.J.  
Deposited on : 2006-12-21  
Resolution : 2.50 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 : 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



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Mol	Chain	Length	Quality of chain
4	E	194	 <p>14% 68% 27% . .</p>
4	L	194	 <p>17% 72% 21% 5% .</p>
5	F	238	 <p>15% 76% 16% 5% .</p>
5	M	238	 <p>8% 84% 14% .</p>

## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 13180 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 HLA CLASS I HISTOCOMPATIBILITY ANTIGEN, A-2 ALPHA CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	275	2238	1395	408	426	9	4	0	0
1	H	275	2238	1395	408	426	9	3	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	167	ALA	TRP	engineered mutation	UNP P01892
H	167	ALA	TRP	engineered mutation	UNP P01892

- Molecule 2 is a protein called BETA-2-MICROGLOBULIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	100	837	533	141	159	4	0	0	0
2	I	100	837	533	141	159	4	0	0	0

- Molecule 3 is a protein called P1049.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
3	C	9	76	56	10	10	0	0	0
3	J	9	76	56	10	10	0	0	0

- Molecule 4 is a protein called TCR ALPHA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	E	194	Total	C	N	O	S	86	0	0
			1521	965	245	302	9			
4	L	194	Total	C	N	O	S	87	0	0
			1521	965	245	302	9			

- Molecule 5 is a protein called TCR BETA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	F	237	Total	C	N	O	S	0	0	0
			1891	1194	331	361	5			
5	M	237	Total	C	N	O	S	3	0	0
			1891	1194	331	361	5			

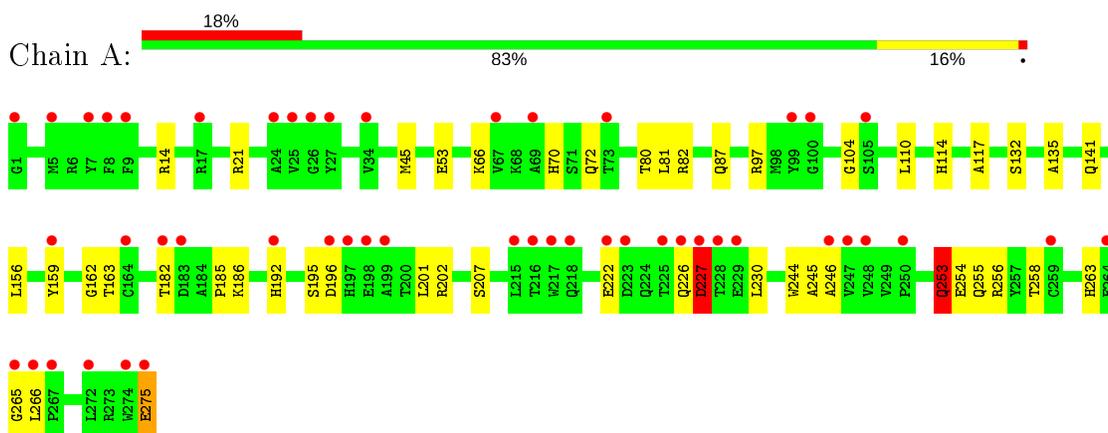
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	9	Total	O	0	0
			9	9		
6	B	1	Total	O	0	0
			1	1		
6	E	10	Total	O	0	0
			10	10		
6	F	5	Total	O	0	0
			5	5		
6	H	9	Total	O	0	0
			9	9		
6	I	6	Total	O	0	0
			6	6		
6	L	9	Total	O	0	0
			9	9		
6	M	5	Total	O	0	0
			5	5		

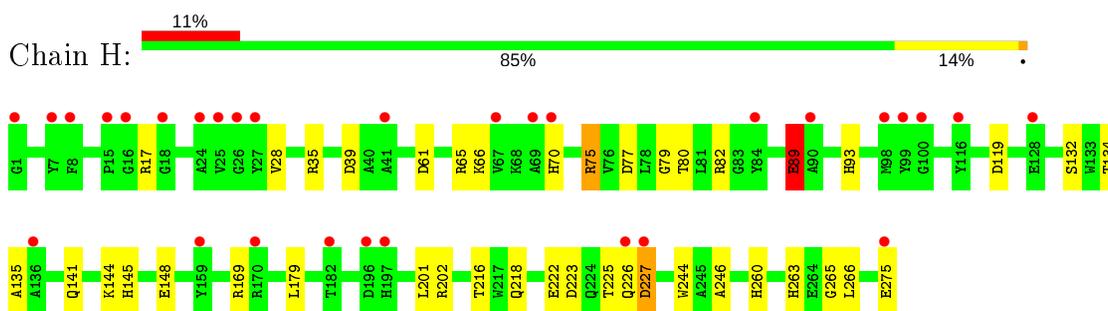
### 3 Residue-property plots [i](#)

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.

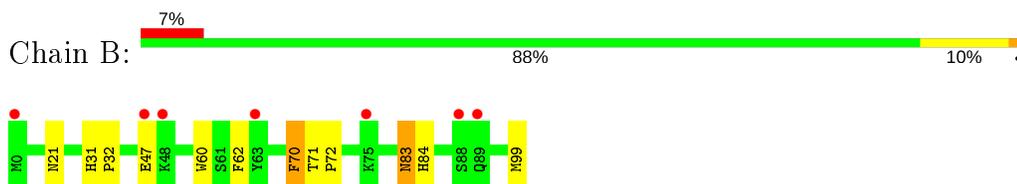
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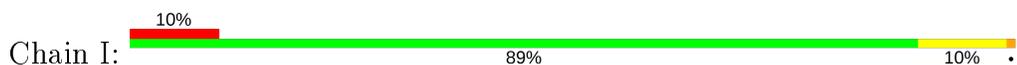
- Molecule 1: HLA CLASS I HISTOCOMPATIBILITY ANTIGEN, A-2 ALPHA CHAIN



- Molecule 2: BETA-2-MICROGLOBULIN

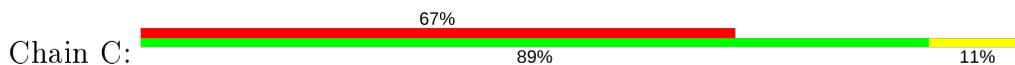


- Molecule 2: BETA-2-MICROGLOBULIN

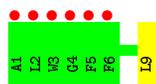
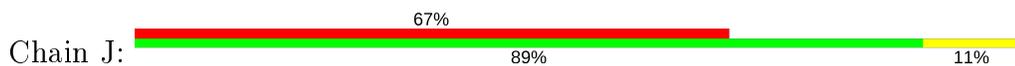




● Molecule 3: P1049



● Molecule 3: P1049



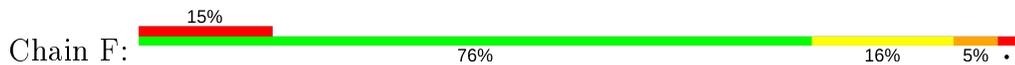
● Molecule 4: TCR ALPHA



● Molecule 4: TCR ALPHA

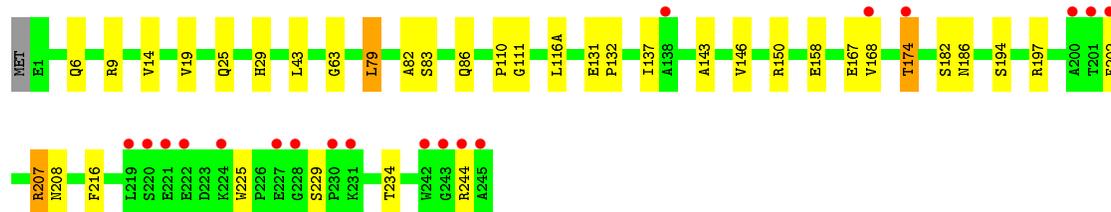


● Molecule 5: TCR BETA



- Molecule 5: TCR BETA

Chain M:  8% 84% 14%



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	94.28Å 84.35Å 122.47Å 90.00° 92.53° 90.00°	Depositor
Resolution (Å)	122.17 – 2.50 40.78 – 2.50	Depositor EDS
% Data completeness (in resolution range)	92.5 (122.17-2.50) 92.5 (40.78-2.50)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.60 (at 2.51Å)	Xtrriage
Refinement program	REFMAC 5.2.0019	Depositor
R, $R_{free}$	0.242 , 0.292 0.269 , 0.298	Depositor DCC
$R_{free}$ test set	3099 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	47.7	Xtrriage
Anisotropy	0.558	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 32.1	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	0.000 for h,-k,-l	Xtrriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	13180	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	50.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 42.51 % 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 2.0198e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

<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

### 5.1 Standard geometry

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	1.07	14/2301 (0.6%)	0.72	7/3121 (0.2%)
1	H	0.76	7/2301 (0.3%)	0.72	1/3121 (0.0%)
2	B	0.73	2/860 (0.2%)	0.65	0/1162
2	I	0.82	3/860 (0.3%)	0.85	1/1162 (0.1%)
3	C	0.81	0/80	0.62	0/108
3	J	0.76	0/80	0.62	0/108
4	E	1.40	24/1555 (1.5%)	1.12	12/2106 (0.6%)
4	L	1.22	14/1557 (0.9%)	1.54	31/2112 (1.5%)
5	F	1.60	29/1947 (1.5%)	1.13	18/2649 (0.7%)
5	M	0.68	3/1947 (0.2%)	0.59	1/2649 (0.0%)
All	All	1.10	96/13488 (0.7%)	0.95	71/18298 (0.4%)

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	1
4	E	0	2
4	L	0	4
5	F	0	4
All	All	0	11

The worst 5 of 96 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	275	GLU	C-O	25.85	1.72	1.23
5	F	62	ASP	C-N	22.68	1.73	1.33
4	E	56	PRO	N-CD	21.81	1.78	1.47
5	F	117	GLU	CD-OE1	20.21	1.47	1.25
5	F	64	GLY	N-CA	19.33	1.75	1.46

The worst 5 of 71 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	F	244	ARG	NE-CZ-NH1	23.41	132.00	120.30
4	L	52	ASP	CA-C-N	-21.03	70.94	117.20
5	F	207	ARG	NE-CZ-NH2	-19.11	110.75	120.30
4	E	134	ARG	NE-CZ-NH1	19.05	129.82	120.30
2	I	99	MET	CA-C-O	18.29	158.52	120.10

There are no chirality outliers.

5 of 11 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	253	GLN	Sidechain
4	E	197	GLU	Sidechain,Mainchain
5	F	153	PHE	Peptide
5	F	204	HIS	Mainchain
5	F	206	PRO	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	2238	0	2091	36	0
1	H	2238	0	2091	25	1
2	B	837	0	803	9	0
2	I	837	0	803	6	0
3	C	76	0	76	2	0
3	J	76	0	76	2	0
4	E	1521	0	1472	32	1
4	L	1521	0	1476	36	0
5	F	1891	0	1793	55	0
5	M	1891	0	1794	21	0
6	A	9	0	0	0	0
6	B	1	0	0	0	0
6	E	10	0	0	0	0
6	F	5	0	0	0	0
6	H	9	0	0	0	0
6	I	6	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	L	9	0	0	1	0
6	M	5	0	0	0	0
All	All	13180	0	12475	202	1

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.

The worst 5 of 202 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)
5:F:134:LYS:CE	5:F:134:LYS:NZ	1.70	1.51
5:F:64:GLY:CA	5:F:64:GLY:N	1.75	1.47
5:F:154:PRO:N	5:F:154:PRO:CD	1.68	1.46
1:A:253:GLN:CD	1:A:253:GLN:NE2	1.69	1.43
5:F:208:ASN:C	5:F:209:HIS:N	1.68	1.41

All (1) 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)
4:E:198:THR:O	1:H:169:ARG:NH2[2_645]	2.14	0.06

## 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	273/275 (99%)	268 (98%)	3 (1%)	2 (1%)	22 39
1	H	273/275 (99%)	262 (96%)	9 (3%)	2 (1%)	22 39
2	B	98/100 (98%)	96 (98%)	2 (2%)	0	100 100
2	I	98/100 (98%)	96 (98%)	2 (2%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	C	7/9 (78%)	6 (86%)	1 (14%)	0	100	100
3	J	7/9 (78%)	6 (86%)	1 (14%)	0	100	100
4	E	189/194 (97%)	165 (87%)	17 (9%)	7 (4%)	3	4
4	L	192/194 (99%)	164 (85%)	17 (9%)	11 (6%)	1	1
5	F	235/238 (99%)	219 (93%)	14 (6%)	2 (1%)	17	31
5	M	235/238 (99%)	227 (97%)	8 (3%)	0	100	100
All	All	1607/1632 (98%)	1509 (94%)	74 (5%)	24 (2%)	10	18

5 of 24 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
4	E	51	THR
4	E	52	ASP
4	E	174	ASP
5	F	154	PRO
4	L	51	THR

### 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	230/230 (100%)	226 (98%)	4 (2%)	60	82
1	H	230/230 (100%)	223 (97%)	7 (3%)	41	68
2	B	95/95 (100%)	93 (98%)	2 (2%)	53	78
2	I	95/95 (100%)	92 (97%)	3 (3%)	39	65
3	C	7/7 (100%)	7 (100%)	0	100	100
3	J	7/7 (100%)	7 (100%)	0	100	100
4	E	177/177 (100%)	169 (96%)	8 (4%)	27	51
4	L	177/177 (100%)	166 (94%)	11 (6%)	18	35
5	F	205/206 (100%)	200 (98%)	5 (2%)	49	74
5	M	205/206 (100%)	199 (97%)	6 (3%)	42	69

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	1428/1430 (100%)	1382 (97%)	46 (3%)	39 65

5 of 46 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	H	39	ASP
1	H	266	LEU
5	M	174	THR
1	H	75	ARG
1	H	132	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 31 such sidechains are listed below:

Mol	Chain	Res	Type
5	F	139	ASN
1	H	93	HIS
5	M	86	GLN
1	H	70	HIS
1	H	141	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
4	E	3
4	L	2
5	F	2

The worst 5 of 7 chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	E	56:PRO	C	57:GLU	N	4.37
1	E	57:GLU	C	58:HIS	N	3.72
1	F	62:ASP	C	64:GLY	N	1.73
1	L	52:ASP	C	53:ASN	N	1.70
1	F	208:ASN	C	209:HIS	N	1.68

## 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	275/275 (100%)	1.09	49 (17%) 1 1	40, 50, 58, 67	1 (0%)
1	H	275/275 (100%)	0.89	30 (10%) 5 5	38, 50, 59, 74	1 (0%)
2	B	100/100 (100%)	0.73	7 (7%) 16 16	42, 49, 59, 73	0
2	I	100/100 (100%)	0.82	10 (10%) 7 6	43, 51, 60, 76	0
3	C	9/9 (100%)	2.00	6 (66%) 0 0	44, 45, 51, 54	0
3	J	9/9 (100%)	2.23	6 (66%) 0 0	48, 50, 54, 57	0
4	E	184/194 (94%)	1.11	27 (14%) 2 2	37, 49, 66, 75	0
4	L	185/194 (95%)	1.08	33 (17%) 1 1	36, 49, 67, 73	2 (1%)
5	F	237/238 (99%)	1.01	35 (14%) 2 2	35, 50, 63, 69	0
5	M	237/238 (99%)	0.73	19 (8%) 12 12	37, 49, 58, 66	1 (0%)
All	All	1611/1632 (98%)	0.96	222 (13%) 2 2	35, 50, 62, 76	5 (0%)

The worst 5 of 222 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
4	E	189	PHE	9.3
5	M	245	ALA	8.7
4	E	173	MET	8.3
5	F	245	ALA	7.8
4	L	156	MET	7.4

### 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.