WORLDWIDE PDB
PROTEIN DATA BANK

Full wwPDB X-ray Structure Validation Report (i)

Feb 1, 2016 – 08:30 AM GMT

PDB ID	:	3EYJ
Title	:	Structure of Influenza Haemagglutinin in complex with an inhibitor of membrane fusion
Authors	:	Russell, R.J.; Kerry, P.S.; Stevens, D.J.; Steinahuer, D.A.; Martin, S.R.; Gamblin, S.J.; Skehel, J.J.
Deposited on	:	2008-10-21
Resolution	:	2.60 Å(reported)

The following versions of software and data (see references (i)) 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 (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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.
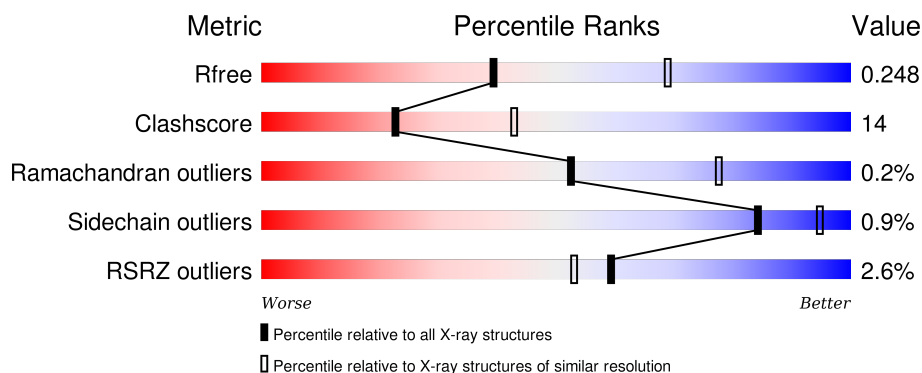
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	91344	2328 (2.60-2.60)
Clashscore	102246	2679 (2.60-2.60)
Ramachandran outliers	100387	2635 (2.60-2.60)
Sidechain outliers	100360	2635 (2.60-2.60)
RSRZ outliers	91569	2334 (2.60-2.60)

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. 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	323	4% — 73% — 23% — ··
2	B	172	84% — 16%

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 4210 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 Hemagglutinin HA1 chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	319	2443	1521	436	475	11	0	0	0

- Molecule 2 is a protein called Hemagglutinin HA2 chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	172	1403	868	251	280	4	0	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
3	A	214	214	214	0	0
3	B	150	150	150	0	0

3 Residue-property plots ⓘ

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: Hemagglutinin HA1 chain
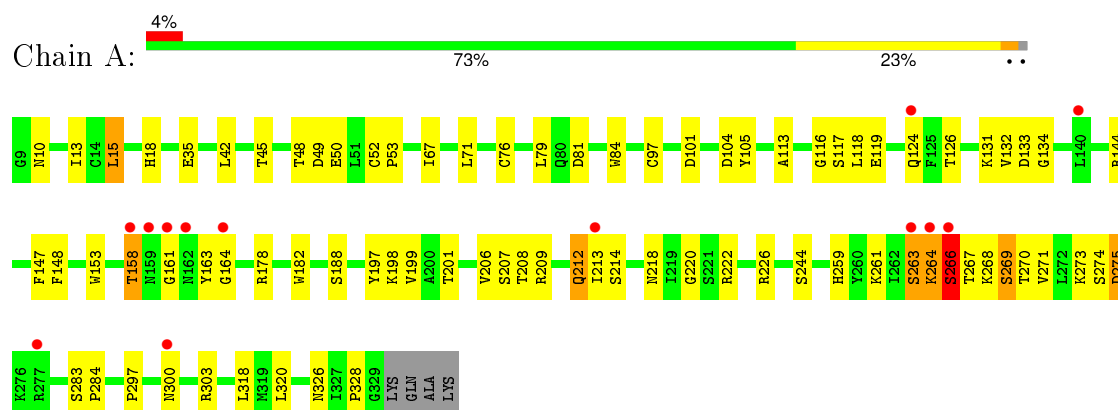
Chain A:

- Molecule 2: Hemagglutinin HA2 chain

Chain B:
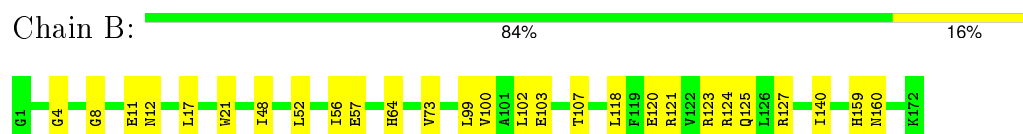
4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 3	Depositor
Cell constants a, b, c, α, β, γ	138.76Å 138.76Å 138.76Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.59 – 2.60 29.58 – 2.60	Depositor EDS
% Data completeness (in resolution range)	51.8 (29.59-2.60) 99.4 (29.58-2.60)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	3.15 (at 2.61Å)	Xtriage
Refinement program	PHENIX (phenix.refine)	Depositor
R, R_{free}	0.193 , 0.244 0.205 , 0.248	Depositor DCC
R_{free} test set	1381 reflections (5.30%)	DCC
Wilson B-factor (Å^2)	33.8	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 44.2	EDS
Estimated twinning fraction	0.029 for l,-k,h	Xtriage
L-test for twinning[2]	$< \lvert L \rvert > = 0.50, < L^2 > = 0.33$	Xtriage
Outliers	0 of 27461 reflections	Xtriage
F_o,F_c correlation	0.93	EDS
Total number of atoms	4210	wwPDB-VP
Average B, all atoms (Å^2)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.40% of the height of the origin peak. No significant pseudotranslation is detected.*

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[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< \lvert L \rvert >$, $< L^2 >$ 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	0.68	2/2493 (0.1%)	0.96	11/3387 (0.3%)
2	B	0.50	0/1426	0.59	0/1919
All	All	0.62	2/3919 (0.1%)	0.84	11/5306 (0.2%)

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

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	158	THR	C-N	12.71	1.63	1.34
1	A	161	GLY	C-N	-6.23	1.19	1.34

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	158	THR	O-C-N	20.79	155.96	122.70
1	A	158	THR	CA-C-N	-15.39	83.33	117.20
1	A	158	THR	C-N-CA	14.34	157.55	121.70
1	A	263	SER	C-N-CA	10.51	147.97	121.70
1	A	264	LYS	O-C-N	10.18	138.98	122.70
1	A	269	SER	N-CA-CB	-8.14	98.29	110.50
1	A	266	SER	C-N-CA	-7.34	103.36	121.70
1	A	15	LEU	CA-CB-CG	-7.01	99.17	115.30
1	A	266	SER	CA-C-O	6.42	133.57	120.10
1	A	269	SER	N-CA-C	6.42	128.33	111.00
1	A	266	SER	CA-C-N	-5.46	105.19	117.20

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	266	SER	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	2443	0	2409	87	2
2	B	1403	0	1324	25	1
3	A	214	0	0	15	0
3	B	150	0	0	5	0
All	All	4210	0	3733	103	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 14.

All (103) 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:131:LYS:NZ	1:A:158:THR:CB	2.07	1.16
1:A:264:LYS:O	1:A:267:THR:CG2	1.96	1.11
1:A:273:LYS:HG2	3:A:537:HOH:O	1.50	1.10
1:A:264:LYS:O	1:A:267:THR:HG23	1.46	1.10
1:A:131:LYS:CE	1:A:158:THR:OG1	2.01	1.08
1:A:45:THR:HB	1:A:300:ASN:HD21	1.18	1.05
1:A:131:LYS:HZ2	1:A:158:THR:CB	1.69	1.02
1:A:263:SER:O	3:A:389:HOH:O	1.78	1.01
1:A:45:THR:HG22	1:A:300:ASN:ND2	1.78	0.98
1:A:266:SER:HB2	3:B:207:HOH:O	1.65	0.95
1:A:131:LYS:HZ1	1:A:158:THR:CB	1.74	0.95
1:A:45:THR:CB	1:A:300:ASN:HD21	1.79	0.95
1:A:268:LYS:HG3	1:A:268:LYS:O	1.66	0.95
1:A:218:ASN:OD1	1:A:222:ARG:NH2	2.00	0.95
1:A:117:SER:OG	1:A:119:GLU:HG2	1.70	0.91
1:A:264:LYS:O	1:A:267:THR:HG22	1.76	0.85

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:45:THR:CG2	1:A:300:ASN:ND2	2.40	0.84
1:A:45:THR:CB	1:A:300:ASN:ND2	2.41	0.83
1:A:131:LYS:NZ	1:A:158:THR:OG1	0.67	0.82
1:A:268:LYS:O	1:A:268:LYS:CG	2.27	0.81
1:A:45:THR:HB	1:A:300:ASN:ND2	1.96	0.81
1:A:104:ASP:HB2	3:A:402:HOH:O	1.80	0.80
1:A:268:LYS:HB3	3:A:514:HOH:O	1.83	0.77
1:A:264:LYS:C	1:A:267:THR:HG23	2.06	0.75
1:A:264:LYS:CE	3:A:442:HOH:O	2.36	0.73
2:B:103:GLU:O	3:B:176:HOH:O	2.07	0.71
1:A:269:SER:O	1:A:270:THR:HB	1.90	0.71
1:A:18:HIS:ND1	2:B:21:TRP:HA	2.07	0.70
1:A:76:CYS:O	1:A:79:LEU:HB2	1.96	0.65
1:A:264:LYS:CG	3:A:442:HOH:O	2.46	0.63
1:A:45:THR:HA	1:A:300:ASN:OD1	1.99	0.63
1:A:131:LYS:HZ1	1:A:158:THR:CG2	2.11	0.62
1:A:45:THR:HG22	1:A:300:ASN:HD22	1.64	0.62
1:A:264:LYS:HE2	3:A:442:HOH:O	1.99	0.62
1:A:303:ARG:HD2	3:B:291:HOH:O	2.00	0.62
1:A:117:SER:HG	1:A:119:GLU:HG2	1.65	0.60
2:B:103:GLU:HA	2:B:103:GLU:OE1	2.02	0.59
1:A:266:SER:CB	3:B:207:HOH:O	2.35	0.59
1:A:199:VAL:HG12	1:A:201:THR:H	1.66	0.58
1:A:263:SER:HB2	1:A:267:THR:HG21	1.87	0.56
2:B:99:LEU:O	2:B:103:GLU:HG2	2.05	0.56
1:A:45:THR:HA	1:A:300:ASN:CG	2.26	0.56
2:B:127:ARG:HG3	2:B:159:HIS:CG	2.40	0.56
1:A:274:SER:OG	1:A:275:ASP:N	2.39	0.55
2:B:56:ILE:HD12	2:B:57:GLU:N	2.22	0.55
1:A:269:SER:O	2:B:64:HIS:HB2	2.06	0.55
2:B:120:GLU:OE1	2:B:123:ARG:NH1	2.41	0.54
1:A:208:THR:HA	1:A:244:SER:O	2.08	0.53
1:A:320:LEU:HD11	2:B:103:GLU:HB3	1.90	0.53
1:A:209:ARG:HD2	3:A:348:HOH:O	2.08	0.53
1:A:18:HIS:N	2:B:17:LEU:HD23	2.24	0.53
1:A:126:THR:HG22	3:A:463:HOH:O	2.08	0.53
1:A:188:SER:HA	1:A:220:GLY:O	2.10	0.51
2:B:4:GLY:O	2:B:8:GLY:HA3	2.10	0.51
1:A:264:LYS:HE3	3:A:442:HOH:O	2.07	0.51
1:A:42:LEU:O	1:A:297:PRO:HD2	2.11	0.51
1:A:264:LYS:O	1:A:266:SER:C	2.49	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:56:ILE:O	2:B:57:GLU:C	2.50	0.50
1:A:147:PHE:O	1:A:148:PHE:C	2.47	0.49
2:B:11:GLU:HG2	2:B:12:ASN:OD1	2.13	0.49
2:B:124:ARG:HD2	3:B:234:HOH:O	2.13	0.48
1:A:318:LEU:HB3	2:B:100:VAL:HG21	1.95	0.48
1:A:84:TRP:CE2	1:A:116:GLY:HA2	2.48	0.48
1:A:264:LYS:HG2	3:A:442:HOH:O	2.11	0.48
1:A:45:THR:HA	1:A:300:ASN:ND2	2.29	0.47
1:A:97:CYS:O	1:A:226:ARG:NH1	2.46	0.47
1:A:50:GLU:HB2	3:A:407:HOH:O	2.15	0.47
1:A:45:THR:CA	1:A:300:ASN:ND2	2.78	0.46
1:A:178:ARG:HD3	1:A:259:HIS:CE1	2.51	0.46
1:A:35:GLU:HG2	1:A:326:ASN:HB3	1.97	0.46
1:A:81:ASP:HA	1:A:119:GLU:HA	1.97	0.46
1:A:67:ILE:HD12	1:A:105:TYR:CZ	2.52	0.45
1:A:52:CYS:HA	1:A:53:PRO:HD3	1.76	0.45
2:B:48:ILE:CD1	2:B:107:THR:HG23	2.47	0.45
1:A:101:ASP:HB3	3:A:516:HOH:O	2.16	0.45
1:A:182:TRP:CE2	1:A:206:VAL:HG21	2.52	0.45
1:A:18:HIS:HB3	2:B:17:LEU:HD23	1.99	0.44
1:A:15:LEU:HD22	2:B:118:LEU:HG	2.00	0.43
2:B:121:ARG:O	2:B:125:GLN:HG3	2.19	0.43
1:A:45:THR:CA	1:A:300:ASN:HD21	2.27	0.43
2:B:73:VAL:O	2:B:73:VAL:HG23	2.18	0.43
1:A:328:PRO:HD2	3:A:265:HOH:O	2.17	0.43
1:A:132:VAL:O	1:A:133:ASP:HB2	2.19	0.43
1:A:212:GLN:HG2	1:A:212:GLN:O	2.14	0.43
1:A:212:GLN:HG3	1:A:213:ILE:N	2.32	0.42
1:A:207:SER:HB3	1:A:212:GLN:HB2	2.01	0.42
1:A:134:GLY:HA3	1:A:153:TRP:HB3	2.01	0.42
1:A:261:LYS:HE2	3:A:393:HOH:O	2.18	0.42
2:B:127:ARG:HG3	2:B:159:HIS:CD2	2.55	0.42
1:A:275:ASP:OD1	1:A:275:ASP:C	2.57	0.42
2:B:99:LEU:O	2:B:103:GLU:CG	2.68	0.42
1:A:67:ILE:HD12	1:A:105:TYR:CE1	2.55	0.41
1:A:283:SER:HA	1:A:284:PRO:HD3	1.93	0.41
1:A:10:ASN:HB3	2:B:140:ILE:O	2.19	0.41
1:A:197:TYR:O	1:A:198:LYS:HB3	2.20	0.41
1:A:131:LYS:NZ	1:A:158:THR:HG1	1.04	0.41
1:A:71:LEU:O	1:A:148:PHE:HB3	2.21	0.41
1:A:13:ILE:HG22	2:B:140:ILE:HD11	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:113:ALA:CB	1:A:271:VAL:HG12	2.51	0.41
1:A:163:TYR:O	1:A:164:GLY:C	2.59	0.41
2:B:52:LEU:HD12	2:B:52:LEU:HA	1.82	0.40
1:A:48:THR:O	1:A:49:ASP:HB3	2.21	0.40
1:A:118:LEU:HA	1:A:118:LEU:HD23	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)
1:A:214:SER:CB	1:A:218:ASN:ND2[9_555]	1.89	0.31
1:A:124:GLN:NE2	2:B:160:ASN:ND2[2_554]	2.06	0.14

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	317/323 (98%)	308 (97%)	8 (2%)	1 (0%)	46	72
2	B	170/172 (99%)	165 (97%)	5 (3%)	0	100	100
All	All	487/495 (98%)	473 (97%)	13 (3%)	1 (0%)	52	77

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	266	SER

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	279/282 (99%)	276 (99%)	3 (1%)	80	93
2	B	147/147 (100%)	146 (99%)	1 (1%)	88	96
All	All	426/429 (99%)	422 (99%)	4 (1%)	84	95

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

Mol	Chain	Res	Type
1	A	144	ARG
1	A	212	GLN
1	A	275	ASP
2	B	102	LEU

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

Mol	Chain	Res	Type
1	A	167	ASN
1	A	248	ASN

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 (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	319/323 (98%)	-0.43	13 (4%)	41	33	19, 31, 53, 69	0
2	B	172/172 (100%)	-0.75	0	100	100	15, 26, 45, 72	0
All	All	491/495 (99%)	-0.54	13 (2%)	59	53	15, 29, 52, 72	0

All (13) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	159	ASN	7.1
1	A	264	LYS	6.1
1	A	158	THR	5.8
1	A	266	SER	5.8
1	A	162	ASN	3.3
1	A	300	ASN	2.8
1	A	164	GLY	2.5
1	A	161	GLY	2.3
1	A	124	GLN	2.3
1	A	213	ILE	2.2
1	A	263	SER	2.1
1	A	277	ARG	2.1
1	A	140	LEU	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.