



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 22, 2022 – 08:10 PM JST

PDB ID : 7FI6
Title : Crystal structure of human MICA mutants in complex with natural killer cell receptor NKG2D
Authors : Cai, W.; Peng, S.; Xu, T.; Tian, Y.; Li, Y.; Liu, J.
Deposited on : 2021-07-30
Resolution : 2.90 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 ⓘ 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 ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.30
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)	:	2.30

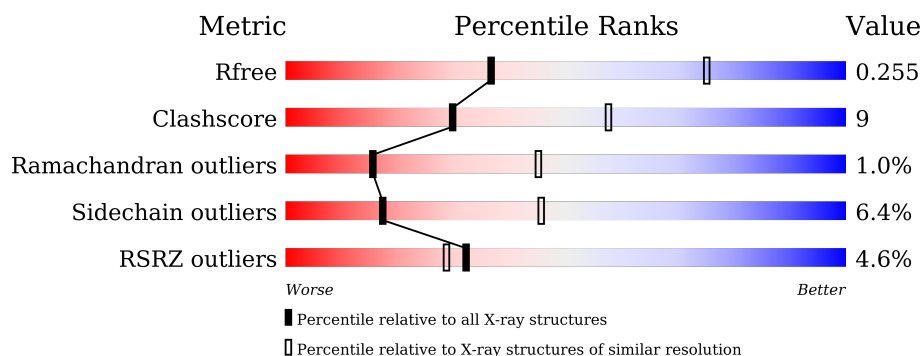
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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	1957 (2.90-2.90)
Clashscore	141614	2172 (2.90-2.90)
Ramachandran outliers	138981	2115 (2.90-2.90)
Sidechain outliers	138945	2117 (2.90-2.90)
RSRZ outliers	127900	1906 (2.90-2.90)

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 $\leq 5\%$. 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	139	<div> <div>6%</div> <div> <div></div> <div>70%</div> <div>19%</div> <div>12%</div> </div> </div>
1	B	139	<div> <div>5%</div> <div> <div></div> <div>68%</div> <div>24%</div> <div>6%</div> </div> </div>
2	C	275	<div> <div>3%</div> <div> <div></div> <div>69%</div> <div>25%</div> <div>.</div> </div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4193 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 NKG2-D type II integral membrane protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	123	Total	C	N	O	S	0	0	0
			996	633	160	191	12			
1	B	131	Total	C	N	O	S	0	0	0
			1060	674	170	204	12			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	78	MET	-	initiating methionine	UNP P26718
A	79	GLU	-	expression tag	UNP P26718
B	78	MET	-	initiating methionine	UNP P26718
B	79	GLU	-	expression tag	UNP P26718

- Molecule 2 is a protein called MHC class I polypeptide-related sequence A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	263	Total	C	N	O	S	0	0	0
			2135	1317	396	408	14			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	0	MET	-	initiating methionine	UNP Q29983
C	161	ARG	HIS	engineered mutation	UNP Q29983
C	177	ILE	VAL	engineered mutation	UNP Q29983

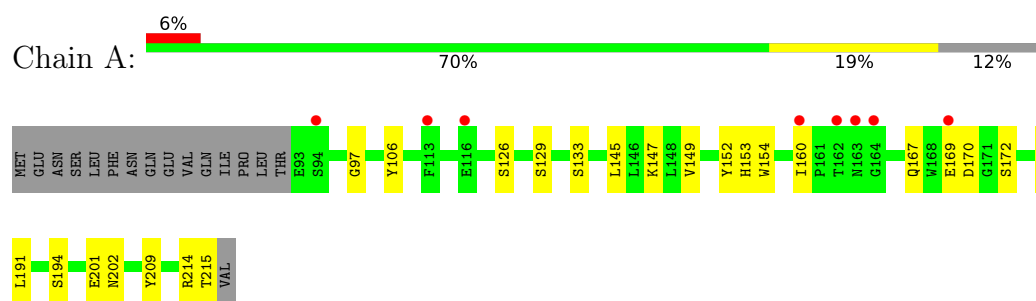
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	C	2	Total	O	0	0
			2	2		

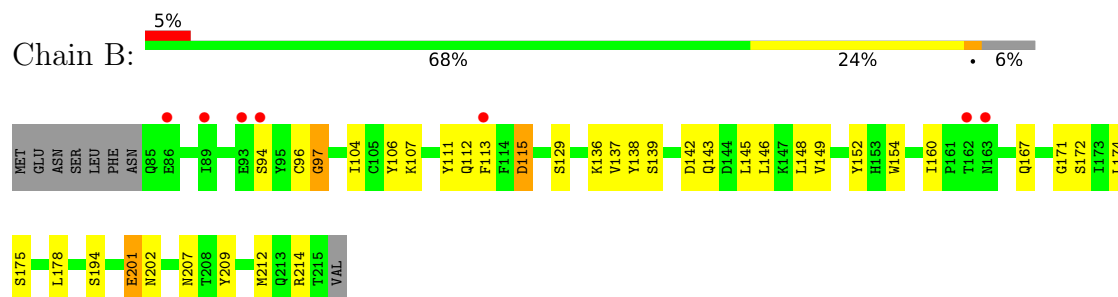
3 Residue-property plots [i](#)

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.

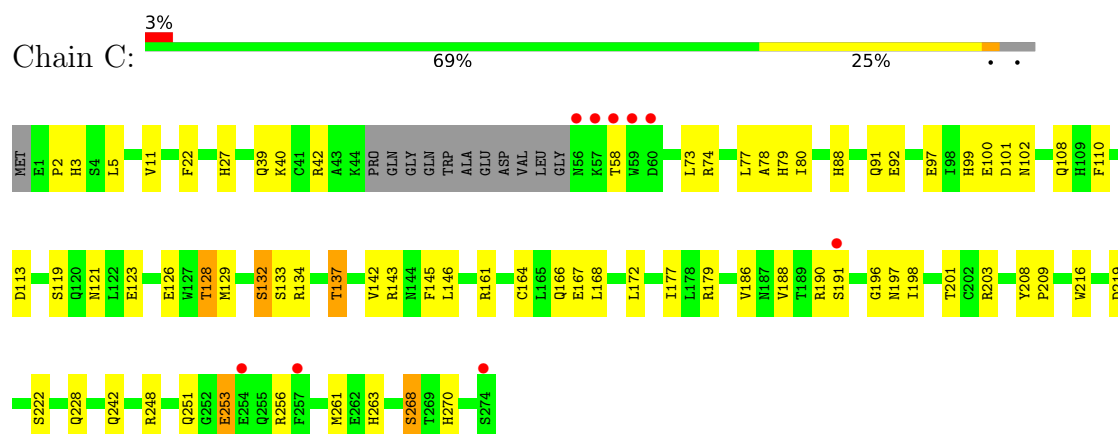
- Molecule 1: NKG2-D type II integral membrane protein



- Molecule 1: NKG2-D type II integral membrane protein



- Molecule 2: MHC class I polypeptide-related sequence A



4 Data and refinement statistics

Property	Value	Source
Space group	I 4 2 2	Depositor
Cell constants a, b, c, α , β , γ	123.60Å 123.60Å 180.77Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	62.91 – 2.90 62.83 – 2.90	Depositor EDS
% Data completeness (in resolution range)	99.9 (62.91-2.90) 99.9 (62.83-2.90)	Depositor EDS
R_{merge}	0.36	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.40 (at 2.91Å)	Xtriage
Refinement program	REFMAC 5.8.0258	Depositor
R, R_{free}	0.193 , 0.262 0.196 , 0.255	Depositor DCC
R_{free} test set	771 reflections (4.85%)	wwPDB-VP
Wilson B-factor (Å ²)	53.4	Xtriage
Anisotropy	0.193	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 39.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.032 for -1/2*h-1/2*k-1/2*l,-1/2*h-1/2*k+1/2*l,-h+k 0.024 for -1/2*h+1/2*k-1/2*l,1/2*h-1/2*k-1/2*l,-h-k	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	4193	wwPDB-VP
Average B, all atoms (Å ²)	58.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.55% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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 [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	0.64	0/1024	0.83	0/1388
1	B	0.64	0/1089	0.78	0/1478
2	C	0.64	0/2183	0.86	0/2951
All	All	0.64	0/4296	0.84	0/5817

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	996	0	933	20	0
1	B	1060	0	1000	29	0
2	C	2135	0	2043	33	0
3	C	2	0	0	0	0
All	All	4193	0	3976	77	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 9.

All (77) 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:145:LEU:HD13	1:B:104:ILE:HD11	1.57	0.86
1:B:160:ILE:HD11	1:B:167:GLN:OE1	1.83	0.78
1:B:136:LYS:O	1:B:212:MET:HE1	1.88	0.74
1:B:137:VAL:HA	1:B:143:GLN:HE22	1.51	0.73
1:B:112:GLN:HE21	1:B:113:PHE:H	1.39	0.69
1:A:191:LEU:CD1	1:A:201:GLU:HG3	2.23	0.68
1:A:185:GLN:HE21	1:A:202:ASN:H	1.42	0.67
2:C:133:SER:O	2:C:137:THR:HG23	1.93	0.67
1:A:191:LEU:HD11	1:A:201:GLU:HG3	1.80	0.64
1:B:106:TYR:CD2	1:B:145:LEU:HD23	2.34	0.63
1:A:145:LEU:HD13	1:B:104:ILE:CD1	2.30	0.60
1:B:104:ILE:HD12	1:B:113:PHE:HE1	1.67	0.60
1:B:201:GLU:HG3	1:B:202:ASN:N	2.16	0.59
2:C:92:GLU:OE2	2:C:108:GLN:NE2	2.34	0.58
1:B:167:GLN:NE2	1:B:171:GLY:O	2.37	0.57
2:C:179:ARG:O	2:C:179:ARG:HG3	2.06	0.56
2:C:191:SER:O	2:C:198:ILE:HG22	2.06	0.55
2:C:5:LEU:HB2	2:C:168:LEU:HD13	1.89	0.55
2:C:216:TRP:H	2:C:228:GLN:NE2	2.04	0.54
1:B:106:TYR:CG	1:B:145:LEU:HD23	2.45	0.52
2:C:2:PRO:HB3	2:C:99:HIS:ND1	2.25	0.52
2:C:91:GLN:O	2:C:110:PHE:HA	2.09	0.52
2:C:203:ARG:NH2	2:C:242:GLN:OE1	2.43	0.52
1:B:138:TYR:H	1:B:143:GLN:HE22	1.57	0.52
1:B:145:LEU:C	1:B:145:LEU:HD12	2.31	0.51
1:A:133:SER:HB2	1:A:169:GLU:OE2	2.11	0.50
1:A:160:ILE:HD12	1:A:167:GLN:OE1	2.12	0.50
1:B:115:ASP:OD1	1:B:115:ASP:N	2.37	0.49
2:C:164:CYS:O	2:C:167:GLU:HB2	2.12	0.49
1:B:138:TYR:H	1:B:143:GLN:NE2	2.09	0.49
1:A:201:GLU:CD	2:C:74:ARG:HH22	2.16	0.49
1:A:152:TYR:CD1	1:A:191:LEU:HD13	2.48	0.48
2:C:253:GLU:O	2:C:256:ARG:HB2	2.14	0.48
1:A:169:GLU:CG	1:A:170:ASP:N	2.76	0.48
1:B:104:ILE:HD12	1:B:113:PHE:CE1	2.49	0.48
1:B:136:LYS:O	1:B:212:MET:CE	2.61	0.48
1:A:106:TYR:CE1	1:A:145:LEU:HD12	2.50	0.47
2:C:209:PRO:O	2:C:263:HIS:HE1	1.99	0.46
2:C:208:TYR:CG	2:C:209:PRO:HA	2.51	0.46
1:B:96:CYS:O	1:B:97:GLY:O	2.34	0.46
1:B:154:TRP:HB2	1:B:209:TYR:HB3	1.98	0.46
1:B:149:VAL:O	1:B:194:SER:HB2	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:169:GLU:HG3	1:A:170:ASP:N	2.31	0.45
2:C:121:ASN:HD21	2:C:123:GLU:HB2	1.81	0.45
2:C:186:VAL:HG22	2:C:261:MET:HE1	1.97	0.45
1:A:154:TRP:HB2	1:A:209:TYR:HB3	1.98	0.45
2:C:172:LEU:HD23	2:C:177:ILE:HD12	1.98	0.45
1:B:152:TYR:CD1	1:B:152:TYR:N	2.84	0.45
1:B:167:GLN:HB2	1:B:172:SER:O	2.16	0.45
1:B:174:LEU:HD23	1:B:175:SER:O	2.17	0.45
2:C:186:VAL:HG22	2:C:261:MET:CE	2.47	0.45
2:C:142:VAL:O	2:C:146:LEU:HG	2.17	0.44
1:A:178:LEU:HD23	1:A:178:LEU:HA	1.89	0.44
2:C:88:HIS:HA	2:C:113:ASP:OD1	2.18	0.44
1:B:107:LYS:HD2	1:B:142:ASP:HA	2.00	0.44
1:A:167:GLN:HB2	1:A:172:SER:O	2.18	0.43
1:A:149:VAL:O	1:A:194:SER:HB2	2.18	0.43
2:C:79:HIS:CD2	2:C:145:PHE:CD1	3.07	0.43
1:A:154:TRP:HA	1:A:154:TRP:CE3	2.53	0.43
2:C:119:SER:HB2	2:C:128:THR:HG23	2.00	0.43
2:C:196:GLY:C	2:C:197:ASN:HD22	2.23	0.43
1:A:147:LYS:O	1:A:194:SER:OG	2.37	0.42
1:B:143:GLN:O	1:B:146:LEU:HB2	2.19	0.42
2:C:101:ASP:O	2:C:102:ASN:HB2	2.19	0.42
2:C:11:VAL:HG12	2:C:22:PHE:HB3	2.00	0.42
2:C:121:ASN:ND2	2:C:123:GLU:HB2	2.35	0.42
2:C:188:VAL:HA	2:C:201:THR:O	2.20	0.42
1:B:160:ILE:HD11	1:B:167:GLN:CD	2.38	0.42
2:C:77:LEU:HA	2:C:80:ILE:HD12	2.02	0.41
1:A:184:MET:HG3	2:C:78:ALA:HB2	2.02	0.41
1:B:175:SER:HB2	1:B:178:LEU:HD12	2.03	0.41
2:C:3:HIS:O	2:C:97:GLU:HA	2.20	0.41
1:A:153:HIS:CE1	1:B:148:LEU:HD11	2.56	0.41
1:B:111:TYR:CE1	1:B:212:MET:HG3	2.56	0.41
2:C:268:SER:OG	2:C:270:HIS:NE2	2.46	0.40
2:C:11:VAL:HG13	2:C:73:LEU:HD23	2.03	0.40
2:C:219:ASP:OD2	2:C:256:ARG:CZ	2.70	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	121/139 (87%)	114 (94%)	6 (5%)	1 (1%)	19	51
1	B	129/139 (93%)	121 (94%)	7 (5%)	1 (1%)	19	51
2	C	259/275 (94%)	242 (93%)	14 (5%)	3 (1%)	13	40
All	All	509/553 (92%)	477 (94%)	27 (5%)	5 (1%)	15	45

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	97	GLY
2	C	100	GLU
1	B	97	GLY
2	C	39	GLN
2	C	132	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	113/129 (88%)	109 (96%)	4 (4%)	36	70
1	B	121/129 (94%)	114 (94%)	7 (6%)	20	50
2	C	237/246 (96%)	218 (92%)	19 (8%)	12	33
All	All	471/504 (94%)	441 (94%)	30 (6%)	17	45

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	126	SER
1	A	129	SER
1	A	214	ARG
1	A	215	THR
1	B	94	SER
1	B	115	ASP
1	B	129	SER
1	B	139	SER
1	B	201	GLU
1	B	207	ASN
1	B	214	ARG
2	C	27	HIS
2	C	40	LYS
2	C	42	ARG
2	C	58	THR
2	C	126	GLU
2	C	128	THR
2	C	129	MET
2	C	132	SER
2	C	134	ARG
2	C	137	THR
2	C	143	ARG
2	C	161	ARG
2	C	166	GLN
2	C	190	ARG
2	C	222	SER
2	C	248	ARG
2	C	251	GLN
2	C	253	GLU
2	C	268	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	185	GLN
1	B	112	GLN
1	B	143	GLN
1	B	167	GLN
2	C	121	ASN
2	C	197	ASN
2	C	228	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 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

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, 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	123/139 (88%)	0.56	8 (6%) 18 14	34, 51, 94, 142	0
1	B	131/139 (94%)	0.52	7 (5%) 26 22	36, 56, 95, 133	0
2	C	263/275 (95%)	0.43	9 (3%) 45 40	30, 52, 98, 133	0
All	All	517/553 (93%)	0.48	24 (4%) 32 29	30, 53, 98, 142	0

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	C	58	THR	4.3
2	C	274	SER	4.0
2	C	191	SER	3.8
1	A	163	ASN	3.3
2	C	57	LYS	3.1
1	A	116	GLU	3.0
1	A	162	THR	3.0
1	A	160	ILE	3.0
1	B	93	GLU	3.0
1	B	162	THR	2.8
2	C	59	TRP	2.7
2	C	254	GLU	2.6
1	A	113	PHE	2.6
1	B	86	GLU	2.5
1	B	163	ASN	2.4
1	B	94	SER	2.4
2	C	56	ASN	2.4
2	C	60	ASP	2.4
1	A	169	GLU	2.4
1	B	113	PHE	2.4
1	B	89	ILE	2.3
1	A	164	GLY	2.3
2	C	257	PHE	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	94	SER	2.2

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.