

# wwPDB X-ray Structure Validation Summary Report (i)

May 22, 2020 – 08:26 pm BST

PDB ID : 3BYT

Title: A complex between a variant of staphylococcal enterotoxin C3 and the variable

domain of the murine T cell receptor beta chain 8.2

Authors : Cho, S.; Eric, J.S.

Deposited on : 2008-01-16

Resolution : 2.30 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
A user guide is available at

https://www.wwpdb.org/validation/2017/XrayValidationReportHelp with specific help available everywhere you see the (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

 $\begin{array}{ccc} & Mol Probity & : & 4.02b\text{-}467 \\ & Xtriage \ (Phenix) & : & 1.13 \end{array}$ 

EDS : 2.11

Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)

Refmac: 5.8.0158

 $\begin{array}{cccc} & CCP4 & : & 7.0.044 \; (Gargrove) \\ Ideal \; geometry \; (proteins) & : & Engh \; \& \; Huber \; (2001) \end{array}$ 

Ideal geometry (DNA, RNA) : Parkinson et al. (1996)

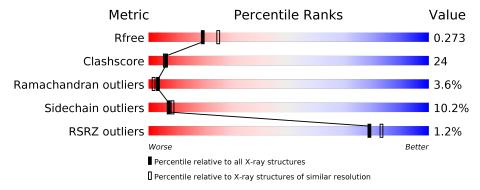
Validation Pipeline (wwPDB-VP) : 2.11

### 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 2.30 Å.

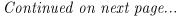
Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	$\begin{array}{c} \text{Whole archive} \\ (\#\text{Entries}) \end{array}$	$\begin{array}{c} {\rm Similar \; resolution} \\ (\#{\rm Entries, \; resolution \; range(\AA)}) \end{array}$
$R_{free}$	130704	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

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 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 <=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	109	48%	40%	10% •			
1	С	109	47%	36%	16% •			
1	Е	109	39%	50%	12%			
1	G	109	48%	40%	10% •			
2	В	239	47%	38%	10% • •			
2	D	239	42%	43%	11% • •			





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Mol	Chain	Length	Oua	lity of chain	
10101	Cham	Dengin	-	inty of chain	
			2%		
2	F	239	44%	41%	11% ••
			<u>%</u>		
2	Н	239	44%	43%	9% • •



## 2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 11054 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 T cell receptor beta chain 8.2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	Λ	109	Total	С	N	О	S	0	0	0
1	A	109	828	513	147	165	3	U	U	
1	С	109	Total	С	N	О	S	0	0	0
1			828	513	147	165	3	U	U	
1	E	109	Total	С	N	О	S	0	0	0
1	نا	109	828	513	147	165	3	0	0	U
1	С	100	Total	С	N	О	S	0	0	0
1	1   G	G 109		513	147	165	3		0	

• Molecule 2 is a protein called Enterotoxin type C-3.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace		
2	В	232	Total	С	N	О	S	0	0	0	
2	Б	232	1900	1205	312	373	10	0	U		
2	D	233	Total	С	N	О	S	0	0	0	
2	D	_ ∠აა	1906	1208	313	375	10	0	U	0	
2	F	234	Total	С	N	О	S	0	0	0	
2	$\Gamma$	234	1917	1215	315	377	10	0	U	$\begin{vmatrix} 0 \end{vmatrix}$	
2	Н	220	Total	С	N	О	S	0	0	0	
	230	1873	1186	307	370	10	U	0	U		

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
В	?	-	VAL	SEE REMARK 999	UNP P0A0L5
В	?	-	GLY	SEE REMARK 999	UNP P0A0L5
В	100B	ALA	-	SEE REMARK 999	UNP P0A0L5
В	101	SER	VAL	SEE REMARK 999	UNP P0A0L5
В	103	TRP	-	SEE REMARK 999	UNP P0A0L5
В	104	HIS	GLY	SEE REMARK 999	UNP P0A0L5
D	?	-	VAL	SEE REMARK 999	UNP P0A0L5
D	?	-	GLY	SEE REMARK 999	UNP P0A0L5



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Chain	Residue	Modelled	Actual	Comment	Reference
D	100B	ALA	-	SEE REMARK 999	UNP P0A0L5
D	101	SER	VAL	SEE REMARK 999	UNP P0A0L5
D	103	TRP	_	SEE REMARK 999	UNP P0A0L5
D	104	HIS	GLY	SEE REMARK 999	UNP P0A0L5
F	?	-	VAL	SEE REMARK 999	UNP P0A0L5
F	?	-	GLY	SEE REMARK 999	UNP P0A0L5
F	100B	ALA	_	SEE REMARK 999	UNP P0A0L5
F	101	SER	VAL	SEE REMARK 999	UNP P0A0L5
F	103	TRP	_	SEE REMARK 999	UNP P0A0L5
F	104	HIS	GLY	SEE REMARK 999	UNP P0A0L5
Н	?	-	VAL	SEE REMARK 999	UNP P0A0L5
Н	?	-	GLY	SEE REMARK 999	UNP P0A0L5
Н	100B	ALA	_	SEE REMARK 999	UNP P0A0L5
Н	101	SER	VAL	SEE REMARK 999	UNP P0A0L5
Н	103	TRP	_	SEE REMARK 999	UNP P0A0L5
Н	104	HIS	GLY	SEE REMARK 999	UNP P0A0L5

#### • Molecule 3 is water.

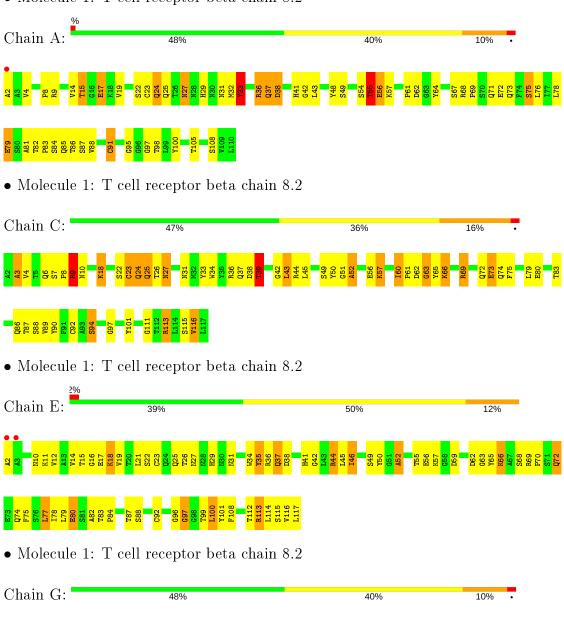
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	11	Total O 11 11	0	0
3	В	21	Total O 21 21	0	0
3	С	14	Total O 14 14	0	0
3	D	27	Total O 27 27	0	0
3	Е	12	Total O 12 12	0	0
3	F	23	Total O 23 23	0	0
3	G	10	Total O 10 10	0	0
3	Н	28	Total O 28 28	0	0



### 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.

• Molecule 1: T cell receptor beta chain 8.2

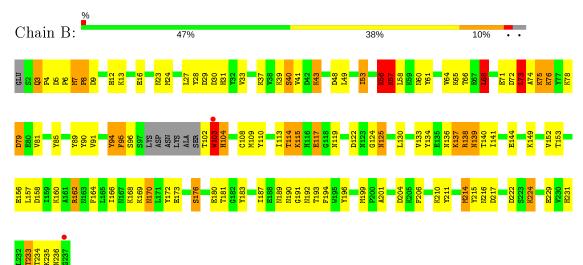




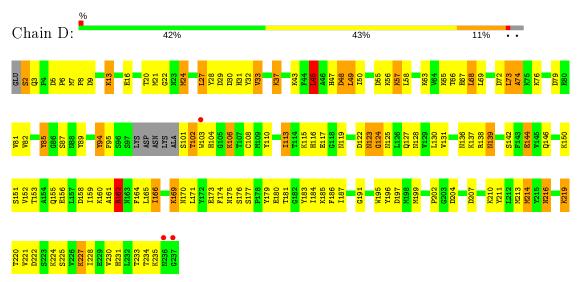




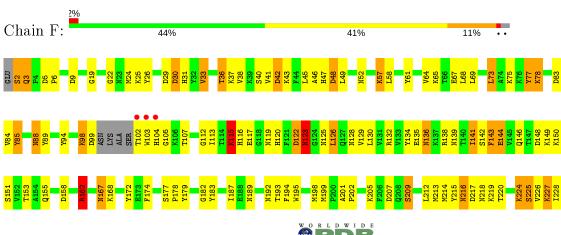
• Molecule 2: Enterotoxin type C-3



• Molecule 2: Enterotoxin type C-3



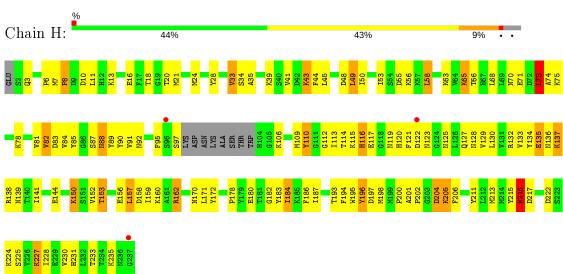
• Molecule 2: Enterotoxin type C-3







• Molecule 2: Enterotoxin type C-3





# 4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1	Depositor
Cell constants	63.16Å 70.64Å 98.84Å	Depositor
a, b, c, $\alpha$ , $\beta$ , $\gamma$	$74.78^{\circ}$ $75.16^{\circ}$ $88.26^{\circ}$	Depositor
Resolution (Å)	30.00 - 2.30	Depositor
Resolution (A)	63.33 - 2.30	EDS
% Data completeness	95.9 (30.00-2.30)	Depositor
(in resolution range)	95.2 (63.33-2.30)	EDS
$R_{merge}$	0.05	Depositor
$R_{sym}$	(Not available)	Depositor
$< I/\sigma(I) > 1$	1.46 (at 2.29Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
P. P.	0.226 , $0.278$	Depositor
$R, R_{free}$	0.222 , $0.273$	DCC
$R_{free}$ test set	3424  reflections  (5.08%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	38.5	Xtriage
Anisotropy	0.308	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$ , $B_{sol}(Å^2)$	0.33 , 51.2	EDS
L-test for twinning <sup>2</sup>	$ < L > = 0.49, < L^2> = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	11054	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 4.45% of the height of the origin peak. No significant pseudotranslation is detected.

<sup>&</sup>lt;sup>2</sup>Theoretical values of <|L|>,  $< L^2>$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.



<sup>&</sup>lt;sup>1</sup>Intensities estimated from amplitudes.

# 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	В	ond lengths	Е	ond angles
WIOI	Chain	RMSZ	# Z  > 5	RMSZ	# Z >5
1	A	1.93	$16/846 \ (1.9\%)$	1.56	11/1145 (1.0%)
1	С	1.96	16/846 (1.9%)	1.66	13/1145 (1.1%)
1	Е	1.84	13/846 (1.5%)	1.51	15/1145~(1.3%)
1	G	1.76	8/846 (0.9%)	1.41	7/1145~(0.6%)
2	В	1.74	22/1943~(1.1%)	1.37	10/2617~(0.4%)
2	D	1.81	34/1949 (1.7%)	1.48	$22/2625 \; (0.8\%)$
2	F	1.73	$26/1960 \ (1.3\%)$	1.41	$17/2639 \ (0.6\%)$
2	Н	1.72	17/1913 (0.9%)	1.38	12/2574~(0.5%)
All	All	1.79	152/11149 (1.4%)	1.45	107/15035~(0.7%)

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 maintain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	С	0	1

The worst 5 of 152 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	$\mathbf{Z}$	$\operatorname{Observed}(\operatorname{\AA})$	$\operatorname{Ideal}(\operatorname{\AA})$
1	E	80	GLU	CG-CD	13.77	1.72	1.51
1	С	52	ALA	CA-CB	12.30	1.78	1.52
1	A	79	GLU	CG-CD	11.22	1.68	1.51
1	A	55	THR	CB-CG2	9.50	1.83	1.52
2	Н	211	TYR	CD2-CE2	-9.36	1.25	1.39

The worst 5 of 107 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\mathbf{Observed}(^o)$	$\operatorname{Ideal}({}^o)$
2	В	162	ARG	NE-CZ-NH1	-13.66	113.47	120.30
1	С	63	GLY	O-C-N	-12.84	102.16	122.70
1	С	9	ARG	NE-CZ-NH1	11.95	126.27	120.30



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Mol	Chain	Res	Type	Atoms	$\mathbf{Z}$	$Observed(^o)$	$\operatorname{Ideal}({}^{o})$
2	D	162	ARG	NE-CZ-NH2	-11.92	114.34	120.30
1	С	9	ARG	NE-CZ-NH2	-10.93	114.84	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	С	63	GLY	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)	$\mathbf{H}(\mathbf{added})$	Clashes	Symm-Clashes
1	A	828	0	788	34	0
1	С	828	0	788	51	0
1	Е	828	0	788	56	0
1	G	828	0	788	48	0
2	В	1900	0	1827	89	0
2	D	1906	0	1832	81	1
2	F	1917	0	1844	94	1
2	Н	1873	0	1804	88	0
3	A	11	0	0	0	0
3	В	21	0	0	1	0
3	С	14	0	0	0	0
3	D	27	0	0	0	0
3	Ε	12	0	0	0	0
3	F	23	0	0	0	0
3	G	10	0	0	1	0
3	Н	28	0	0	0	0
All	All	11054	0	10459	521	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 24.

The worst 5 of 521 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.



Atom-1	Atom-2	$egin{array}{ll}  ext{Interatomic} \  ext{distance } ( ext{Å}) \end{array}$	$egin{aligned}  ext{Clash} \  ext{overlap } ( ext{Å}) \end{aligned}$
2:D:27:LEU:CD1	2:D:27:LEU:CG	1.75	1.60
1:C:52:ALA:CA	1:C:52:ALA:CB	1.78	1.60
1:C:39:THR:CG2	1:C:39:THR:CB	1.75	1.58
1:A:55:THR:CB	1:A:55:THR:CG2	1.83	1.53
2:F:115:LYS:HE3	2:F:115:LYS:N	1.66	1.10

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	$egin{array}{c}  ext{Interatomic} \  ext{distance} \ ( ext{Å}) \end{array}$	$egin{array}{c}  ext{Clash} \  ext{overlap } ( ext{Å}) \end{array}$
2:D:56:LYS:NZ	2:F:123:ASN:OD1[1_554]	2.08	0.12

## 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	Perce	ntiles
1	A	107/109~(98%)	98 (92%)	7 (6%)	2 (2%)	8	7
1	С	107/109 (98%)	97 (91%)	6 (6%)	4 (4%)	3	2
1	E	107/109 (98%)	90 (84%)	14 (13%)	3 (3%)	5	3
1	G	107/109 (98%)	93 (87%)	9 (8%)	5 (5%)	2	1
2	В	$228/239 \ (95\%)$	183 (80%)	$34\ (15\%)$	11 (5%)	2	1
2	D	$229/239 \; (96\%)$	195 (85%)	30 (13%)	4 (2%)	9	8
2	F	230/239~(96%)	183 (80%)	35 (15%)	12 (5%)	2	1
2	Н	$226/239 \; (95\%)$	197 (87%)	22 (10%)	7 (3%)	4	2
All	All	1341/1392 (96%)	1136 (85%)	157 (12%)	48 (4%)	3	2

5 of 48 Ramachandran outliers are listed below:

$\mathbf{Mol}$	Chain	${f Res}$	Type
2	В	57	LYS



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Mol	Chain	Res	Type
2	В	104	HIS
2	В	125	ASN
2	F	57	LYS
2	F	236	ASN

#### 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	88/88 (100%)	76 (86%)	12 (14%)	3 3
1	С	88/88 (100%)	78 (89%)	10 (11%)	5 6
1	E	88/88 (100%)	78 (89%)	10 (11%)	5 6
1	G	88/88 (100%)	77 (88%)	11 (12%)	4 5
2	В	$214/220 \ (97\%)$	193 (90%)	21 (10%)	8 9
2	D	215/220 (98%)	195 (91%)	20 (9%)	9 10
2	F	216/220 (98%)	192 (89%)	24 (11%)	6 7
2	Н	211/220 (96%)	196 (93%)	15 (7%)	14 19
All	All	$1208/1232 \ (98\%)$	1085 (90%)	123 (10%)	7 8

5 of 123 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	D	138	ARG
1	E	46	ILE
2	Н	68	LEU
2	D	144	GLU
2	D	234	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 53 such sidechains are listed below:

Mol	Chain	Res	Type
2	D	136	ASN



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Mol	Chain	Res	Type
1	E	30	ASN
2	Н	92	ASN
2	D	155	GLN
2	D	175	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,  $95^{th}$  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>	$\#\mathrm{RSRZ}{>}2$	$\mathbf{OWAB}(\mathrm{\AA}^2)$	Q < 0.9
1	A	$109/109 \; (100\%)$	-0.19	1 (0%) 84 88	24, 39, 60, 68	0
1	С	109/109 (100%)	-0.26	0 100 100	21, 34, 55, 64	0
1	E	109/109 (100%)	-0.22	2 (1%) 68 74	25, 39, 60, 71	0
1	G	109/109 (100%)	-0.20	0 100 100	26, 46, 67, 76	0
2	В	$232/239 \ (97\%)$	-0.21	2 (0%) 84 88	23, 44, 75, 88	0
2	D	233/239 (97%)	-0.29	3 (1%) 77 81	18, 38, 62, 84	0
2	F	$234/239 \ (97\%)$	-0.09	5 (2%) 63 70	26, 46, 76, 105	0
2	Н	230/239~(96%)	-0.25	3 (1%) 77 81	25, 43, 72, 90	0
All	All	$1365/1392 \ (98\%)$	-0.21	16 (1%) 79 83	18, 42, 69, 105	0

The worst 5 of 16 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	2	ALA	4.1
2	D	237	GLY	4.1
2	D	236	ASN	3.9
2	Н	237	GLY	3.6
2	F	104	HIS	3.6

### 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.

