



wwPDB X-ray Structure Validation Summary Report ⓘ

Jan 31, 2016 – 10:14 PM GMT

PDB ID : 1SLQ
Title : Crystal structure of the trimeric state of the rhesus rotavirus VP4 membrane interaction domain, VP5CT
Authors : Dormitzer, P.R.; Nason, E.B.; Prasad, B.V.V.; Harrison, S.C.
Deposited on : 2004-03-06
Resolution : 3.20 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

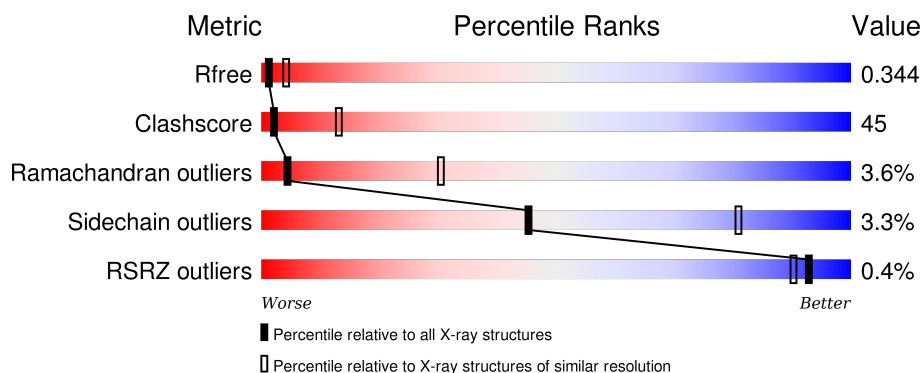
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.20 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.



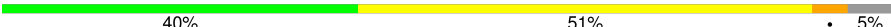
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	91344	1124 (3.24-3.16)
Clashscore	102246	1024 (3.22-3.18)
Ramachandran outliers	100387	1004 (3.22-3.18)
Sidechain outliers	100360	1003 (3.22-3.18)
RSRZ outliers	91569	1129 (3.24-3.16)

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 $\leq 5\%$. 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	278	<div> <div>%</div> <div> <div></div> <div>37%</div> <div>58%</div> <div>• •</div> </div> </div>
1	B	278	<div> <div>35%</div> <div>57%</div> <div>5%</div> <div>•</div> </div>
1	C	278	<div> <div>41%</div> <div>51%</div> <div>• •</div> </div>
1	D	278	<div> <div>%</div> <div>37%</div> <div>55%</div> <div>• •</div> </div>
1	E	278	<div> <div>39%</div> <div>50%</div> <div>6%</div> <div>•</div> </div>

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Mol	Chain	Length	Quality of chain
1	F	278	 A horizontal bar chart showing the quality of the chain. The bar is divided into three segments: green (40%), yellow (51%), and orange (5%). The green segment is labeled '40%', the yellow segment is labeled '51%', and the orange segment is labeled '5%'.

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 12772 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.

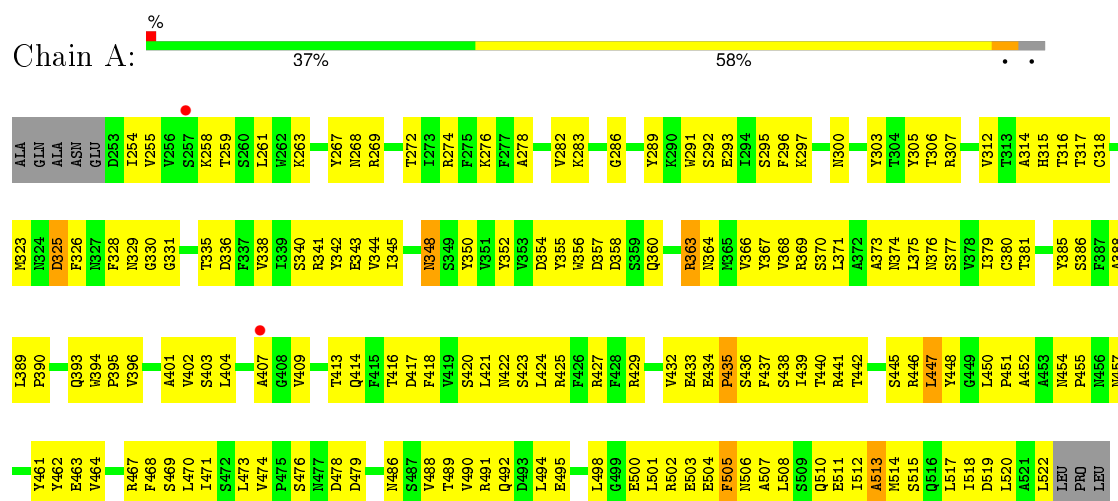
- Molecule 1 is a protein called VP4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	270	Total	C	N	O	S	13	0	0
			2148	1362	361	418	7			
1	B	269	Total	C	N	O	S	0	0	0
			2140	1358	360	415	7			
1	C	266	Total	C	N	O	S	0	0	0
			2120	1342	357	414	7			
1	D	267	Total	C	N	O	S	26	0	0
			2127	1349	358	413	7			
1	E	267	Total	C	N	O	S	0	0	0
			2126	1348	358	413	7			
1	F	265	Total	C	N	O	S	98	0	0
			2111	1337	356	411	7			

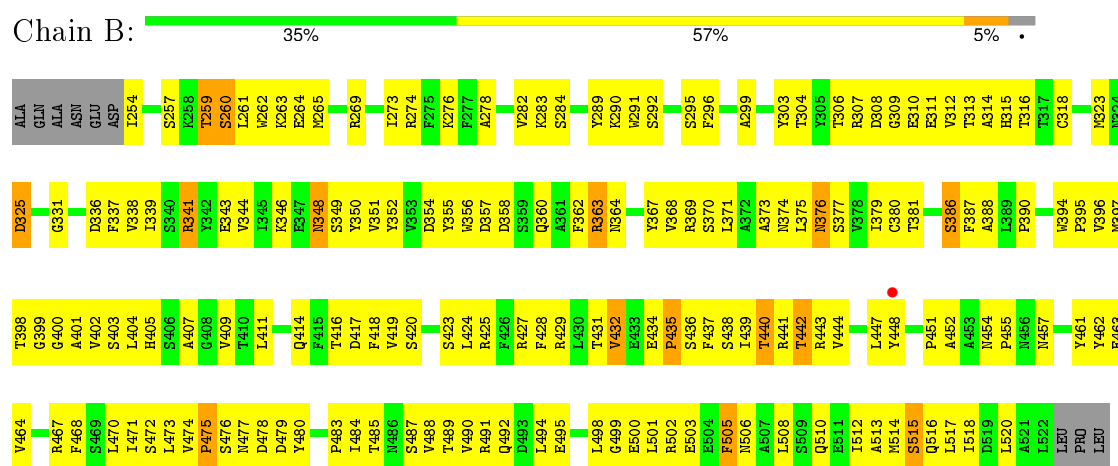
3 Residue-property plots

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 ($\text{RSRZ} > 2$). 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.

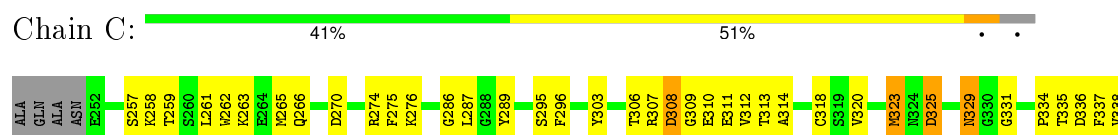
• Molecule 1: VP4

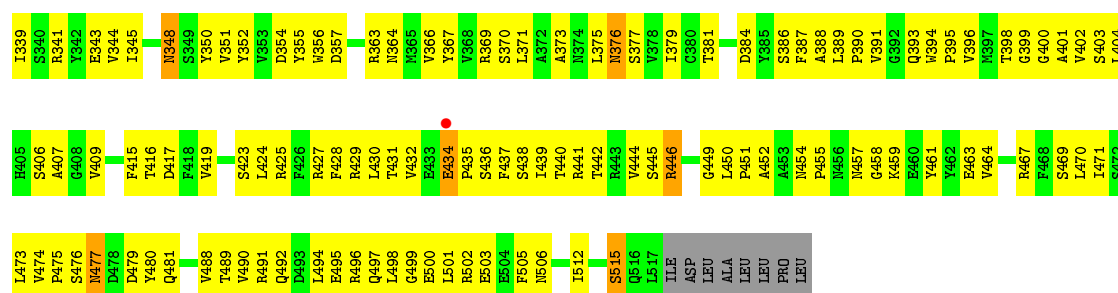


• Molecule 1: VP4

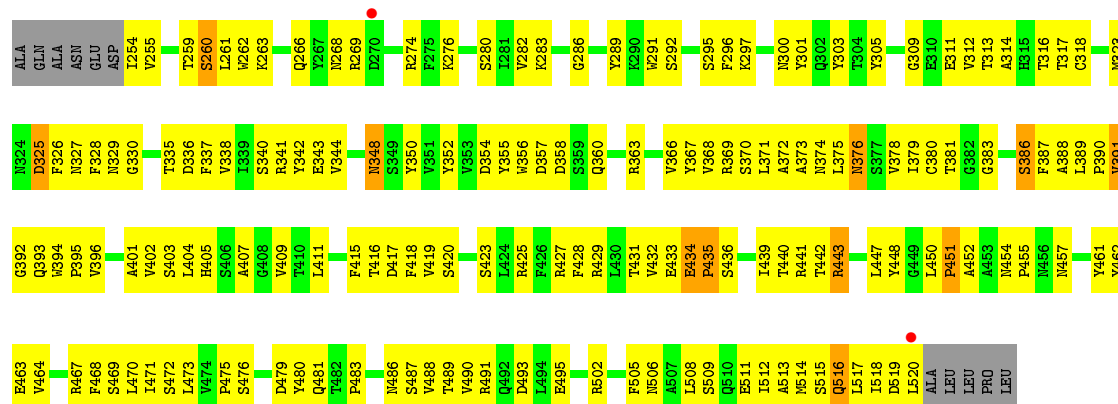


• Molecule 1: VP4

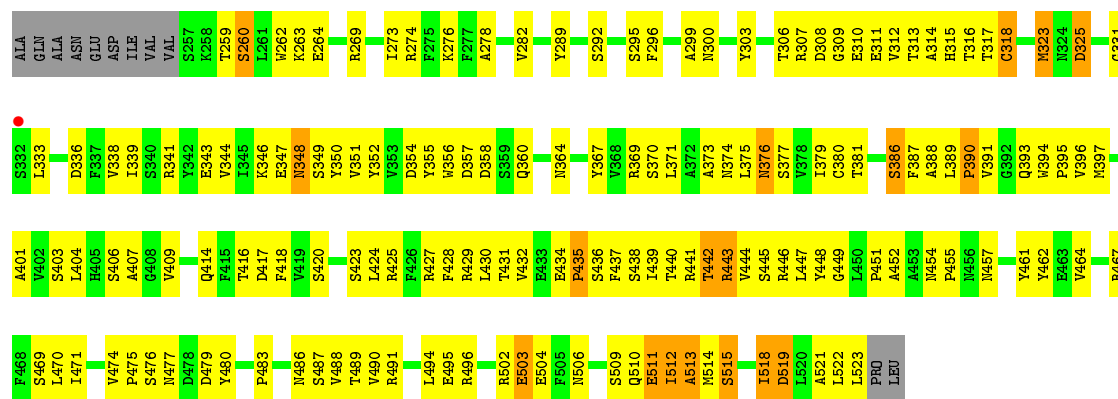




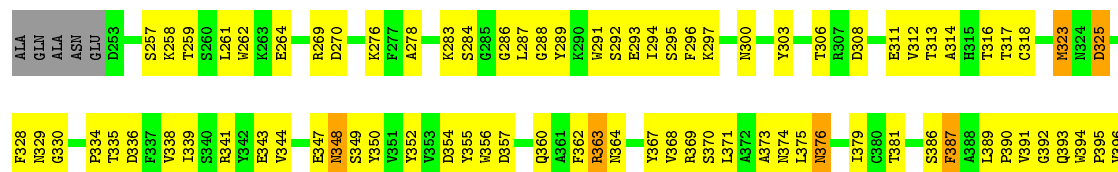
• Molecule 1: VP4

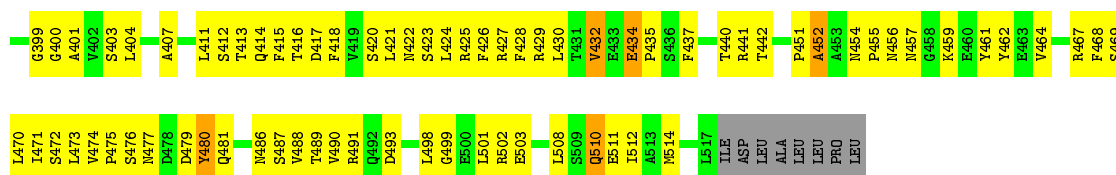


• Molecule 1: VP4



• Molecule 1: VP4





4 Data and refinement statistics

Property	Value	Source
Space group	P 4 ₂ 2 ₂	Depositor
Cell constants a, b, c, α , β , γ	244.84Å 244.84Å 104.78Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 3.20 19.98 – 3.20	Depositor EDS
% Data completeness (in resolution range)	99.8 (20.00-3.20) 99.8 (19.98-3.20)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	12.30 (at 3.22Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.308 , 0.338 0.313 , 0.344	Depositor DCC
R_{free} test set	2625 reflections (4.98%)	DCC
Wilson B-factor (Å ²)	92.3	Xtriage
Anisotropy	0.019	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.24 , 78.9	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.26$	Xtriage
Outliers	1 of 52685 reflections (0.002%)	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	12772	wwPDB-VP
Average B, all atoms (Å ²)	91.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 70.67 % 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 3.0016e-06. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² 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	0.55	0/2196	0.78	0/2983
1	B	0.55	0/2188	0.75	0/2972
1	C	0.55	0/2168	0.77	0/2944
1	D	0.53	0/2175	0.78	0/2954
1	E	0.54	0/2174	0.76	0/2952
1	F	0.54	0/2159	0.79	1/2932 (0.0%)
All	All	0.54	0/13060	0.77	1/17737 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	434	GLU	N-CA-C	5.42	125.65	111.00

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	2148	0	2077	224	0
1	B	2140	0	2073	213	0
1	C	2120	0	2041	213	0
1	D	2127	0	2057	225	0
1	E	2126	0	2055	205	0
1	F	2111	0	2035	194	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	12772	0	12338	1118	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 45.

The worst 5 of 1118 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:D:289:TYR:CD2	1:D:335:THR:HA	1.78	1.18
1:D:286:GLY:O	1:E:510:GLN:HA	1.42	1.18
1:A:439:ILE:HB	1:A:442:THR:HG21	1.26	1.18
1:D:289:TYR:HD2	1:D:335:THR:HA	1.01	1.15
1:A:289:TYR:CD2	1:A:335:THR:HA	1.84	1.13

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	268/278 (96%)	223 (83%)	38 (14%)	7 (3%)	7	40
1	B	267/278 (96%)	222 (83%)	32 (12%)	13 (5%)	3	22
1	C	264/278 (95%)	225 (85%)	32 (12%)	7 (3%)	6	39
1	D	265/278 (95%)	216 (82%)	38 (14%)	11 (4%)	3	26
1	E	265/278 (95%)	219 (83%)	35 (13%)	11 (4%)	3	26
1	F	263/278 (95%)	227 (86%)	28 (11%)	8 (3%)	5	35
All	All	1592/1668 (95%)	1332 (84%)	203 (13%)	57 (4%)	4	30

5 of 57 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	435	PRO
1	A	513	ALA
1	B	440	THR
1	B	475	PRO
1	D	260	SER

5.3.2 Protein sidechains ⓘ

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	238/244 (98%)	231 (97%)	7 (3%)	50	83
1	B	237/244 (97%)	230 (97%)	7 (3%)	48	82
1	C	235/244 (96%)	229 (97%)	6 (3%)	54	85
1	D	236/244 (97%)	228 (97%)	8 (3%)	44	80
1	E	235/244 (96%)	224 (95%)	11 (5%)	32	73
1	F	234/244 (96%)	227 (97%)	7 (3%)	48	82
All	All	1415/1464 (97%)	1369 (97%)	46 (3%)	45	81

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

Mol	Chain	Res	Type
1	D	323	MET
1	D	443	ARG
1	F	348	ASN
1	D	325	ASP
1	D	386	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 63 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	374	ASN
1	D	324	ASN
1	F	329	ASN
1	C	376	ASN

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Mol	Chain	Res	Type
1	C	422	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	268/278 (96%)	-0.27	2 (0%) 89 83	35, 86, 141, 172	0
1	B	269/278 (96%)	-0.39	1 (0%) 93 90	28, 87, 135, 164	0
1	C	266/278 (95%)	-0.33	1 (0%) 93 90	38, 88, 146, 189	0
1	D	264/278 (94%)	-0.29	2 (0%) 87 80	34, 90, 142, 177	0
1	E	267/278 (96%)	-0.32	1 (0%) 93 90	44, 87, 142, 175	0
1	F	253/278 (91%)	-0.36	0 100 100	33, 85, 133, 167	0
All	All	1587/1668 (95%)	-0.33	7 (0%) 93 90	28, 87, 141, 189	0

The worst 5 of 7 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	434	GLU	2.7
1	D	270	ASP	2.6
1	D	520	LEU	2.4
1	A	407	ALA	2.4
1	E	332	SER	2.3

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.