



wwPDB X-ray Structure Validation Summary Report ⓘ

May 22, 2020 – 10:19 pm BST

PDB ID : 3DEX
Title : Crystal structure of SAV_2001 protein from Streptomyces avermitilis, Northeast Structural Genomics Consortium Target SvR107.
Authors : Forouhar, F.; Neely, H.; Seetharaman, J.; Janjua, H.; Fang, Y.; Xiao, R.; Cunningham, K.; Ma, L.-C.; Owen, L.A.; Chen, C.X.; Acton, T.B.; Montelione, G.T.; Hunt, J.F.; Tong, L.; Northeast Structural Genomics Consortium (NESG)
Deposited on : 2008-06-10
Resolution : 2.70 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 ⓘ symbol.

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

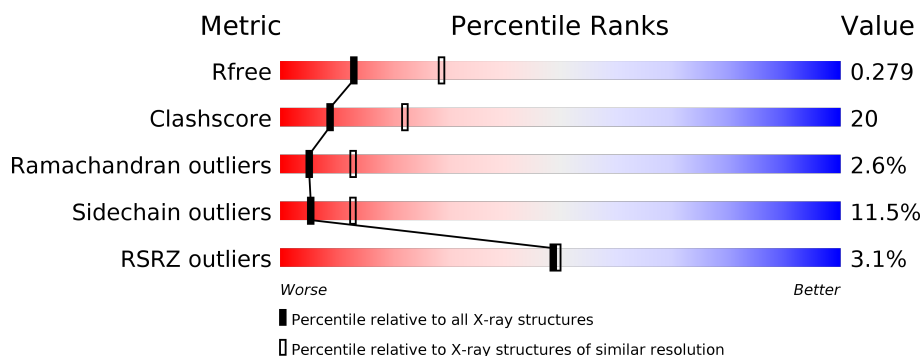
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	2808 (2.70-2.70)
Clashscore	141614	3122 (2.70-2.70)
Ramachandran outliers	138981	3069 (2.70-2.70)
Sidechain outliers	138945	3069 (2.70-2.70)
RSRZ outliers	127900	2737 (2.70-2.70)

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 $\leq 5\%$. 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	107	<div> <div>4%</div> <div> <div></div> <div>37%</div> <div>34%</div> <div>•</div> <div>26%</div> </div> </div>
1	B	107	<div> <div>2%</div> <div> <div></div> <div>37%</div> <div>34%</div> <div>•</div> <div>25%</div> </div> </div>
1	C	107	<div> <div>0%</div> <div> <div></div> <div>36%</div> <div>32%</div> <div>5%</div> <div>28%</div> </div> </div>
1	D	107	<div> <div>3%</div> <div> <div></div> <div>36%</div> <div>35%</div> <div>6%</div> <div>24%</div> </div> </div>
1	E	107	<div> <div>3%</div> <div> <div></div> <div>35%</div> <div>34%</div> <div>6%</div> <div>26%</div> </div> </div>
1	F	107	<div> <div>3%</div> <div> <div></div> <div>37%</div> <div>32%</div> <div>5%</div> <div>26%</div> </div> </div>

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Mol	Chain	Length	Quality of chain
1	G	107	<div><div><div></div><div></div><div></div><div></div></div><div><div>2%</div><div>37%</div><div>32%</div><div>•</div><div>28%</div></div></div>
1	H	107	<div><div><div></div><div></div><div></div><div></div></div><div><div>2%</div><div>37%</div><div>34%</div><div>7%</div><div>22%</div></div></div>

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 5209 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 SAV_2001.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	79	Total	C	N	O	S	0	0	0
			648	411	119	116	2			
1	B	80	Total	C	N	O	S	0	0	0
			656	417	119	118	2			
1	C	77	Total	C	N	O	S	0	0	0
			633	402	115	114	2			
1	D	81	Total	C	N	O	S	0	0	0
			663	421	120	120	2			
1	E	79	Total	C	N	O	S	0	0	0
			648	411	119	116	2			
1	F	79	Total	C	N	O	S	0	0	0
			648	411	119	116	2			
1	G	77	Total	C	N	O	S	0	0	0
			633	402	115	114	2			
1	H	83	Total	C	N	O	S	0	0	0
			680	432	124	122	2			

There are 64 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	100	LEU	-	expression tag	UNP Q82LK9
A	101	GLU	-	expression tag	UNP Q82LK9
A	102	HIS	-	expression tag	UNP Q82LK9
A	103	HIS	-	expression tag	UNP Q82LK9
A	104	HIS	-	expression tag	UNP Q82LK9
A	105	HIS	-	expression tag	UNP Q82LK9
A	106	HIS	-	expression tag	UNP Q82LK9
A	107	HIS	-	expression tag	UNP Q82LK9
B	100	LEU	-	expression tag	UNP Q82LK9
B	101	GLU	-	expression tag	UNP Q82LK9
B	102	HIS	-	expression tag	UNP Q82LK9
B	103	HIS	-	expression tag	UNP Q82LK9
B	104	HIS	-	expression tag	UNP Q82LK9

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Chain	Residue	Modelled	Actual	Comment	Reference
B	105	HIS	-	expression tag	UNP Q82LK9
B	106	HIS	-	expression tag	UNP Q82LK9
B	107	HIS	-	expression tag	UNP Q82LK9
C	100	LEU	-	expression tag	UNP Q82LK9
C	101	GLU	-	expression tag	UNP Q82LK9
C	102	HIS	-	expression tag	UNP Q82LK9
C	103	HIS	-	expression tag	UNP Q82LK9
C	104	HIS	-	expression tag	UNP Q82LK9
C	105	HIS	-	expression tag	UNP Q82LK9
C	106	HIS	-	expression tag	UNP Q82LK9
C	107	HIS	-	expression tag	UNP Q82LK9
D	100	LEU	-	expression tag	UNP Q82LK9
D	101	GLU	-	expression tag	UNP Q82LK9
D	102	HIS	-	expression tag	UNP Q82LK9
D	103	HIS	-	expression tag	UNP Q82LK9
D	104	HIS	-	expression tag	UNP Q82LK9
D	105	HIS	-	expression tag	UNP Q82LK9
D	106	HIS	-	expression tag	UNP Q82LK9
D	107	HIS	-	expression tag	UNP Q82LK9
E	100	LEU	-	expression tag	UNP Q82LK9
E	101	GLU	-	expression tag	UNP Q82LK9
E	102	HIS	-	expression tag	UNP Q82LK9
E	103	HIS	-	expression tag	UNP Q82LK9
E	104	HIS	-	expression tag	UNP Q82LK9
E	105	HIS	-	expression tag	UNP Q82LK9
E	106	HIS	-	expression tag	UNP Q82LK9
E	107	HIS	-	expression tag	UNP Q82LK9
F	100	LEU	-	expression tag	UNP Q82LK9
F	101	GLU	-	expression tag	UNP Q82LK9
F	102	HIS	-	expression tag	UNP Q82LK9
F	103	HIS	-	expression tag	UNP Q82LK9
F	104	HIS	-	expression tag	UNP Q82LK9
F	105	HIS	-	expression tag	UNP Q82LK9
F	106	HIS	-	expression tag	UNP Q82LK9
F	107	HIS	-	expression tag	UNP Q82LK9
G	100	LEU	-	expression tag	UNP Q82LK9
G	101	GLU	-	expression tag	UNP Q82LK9
G	102	HIS	-	expression tag	UNP Q82LK9
G	103	HIS	-	expression tag	UNP Q82LK9
G	104	HIS	-	expression tag	UNP Q82LK9
G	105	HIS	-	expression tag	UNP Q82LK9
G	106	HIS	-	expression tag	UNP Q82LK9

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Chain	Residue	Modelled	Actual	Comment	Reference
G	107	HIS	-	expression tag	UNP Q82LK9
H	100	LEU	-	expression tag	UNP Q82LK9
H	101	GLU	-	expression tag	UNP Q82LK9
H	102	HIS	-	expression tag	UNP Q82LK9
H	103	HIS	-	expression tag	UNP Q82LK9
H	104	HIS	-	expression tag	UNP Q82LK9
H	105	HIS	-	expression tag	UNP Q82LK9
H	106	HIS	-	expression tag	UNP Q82LK9
H	107	HIS	-	expression tag	UNP Q82LK9

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

Chain A:

4% 37% 34% 26%

MET VAL GLY ALA PRO ARG THR THR PRO H11 T12 H13 Q16 I17 E18 T21 Q22 L26 P27 W31 A33 Q34 E35 L36 L37 T38 E41 T42 E43 L44 T45 E46 L47 K50 P51 G52 T53 F57 V58 D62 D63 V66 R69 E71 P75 E76

Chain B:

2% 37% 34% 25%

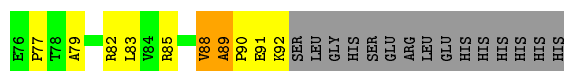
Residue	Category
VAL	Green
GLY	Green
ALA	Green
PRO	Green
ARG	Green
MET	Green
THR	Green
PRO	Green
HIS	Green
THR	Green
H13	Green
R14	Green
V15	Green
Q16	Green
T21	Green
Q22	Green
L26	Green
P27	Green
W31	Green
L32	Green
A33	Green
Q34	Green
E35	Green
L36	Green
L37	Green
T38	Green
E41	Yellow
T42	Yellow
E43	Yellow
L44	Yellow
T45	Yellow
E46	Yellow
L47	Yellow
K50	Yellow
F51	Yellow
G52	Yellow
T53	Yellow
G54	Yellow
G55	Yellow
V56	Yellow
F57	Yellow
V58	Yellow
D62	Yellow
D63	Yellow
V66	Yellow
R69	Yellow
R70	Yellow
E71	Yellow
D75	Yellow

Chain C:

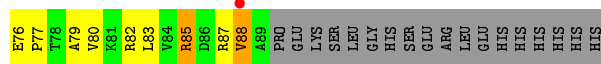
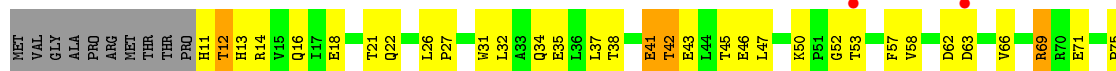
Amino Acid	Percentage
MET	36%
VAL	32%
GLY	5%
ALA	28%
PRO	
ARG	
MET	
THR	
THR	
PRO	
HIS	
T12	
H13	
R14	
V15	
Q16	
T21	
Q22	
L26	
P27	
W31	
L32	
A33	
Q34	
E35	
L36	
L37	
L38	
E41	
T42	
L47	
K50	
P51	
G52	
T53	
G54	
F57	
V58	
D62	
D63	
E64	
V65	
V66	
R69	
R70	
E71	
P75	
E76	
P77	
T78	

Chain D:

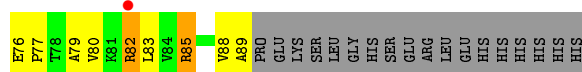
Amino Acid	Count
MET	1
VAL	1
GLY	1
ALA	1
PRO	1
ARG	1
MET	1
THR	1
THR	1
HIS	1
T12	1
H13	1
R14	1
V15	1
Q16	1
T21	1
Q22	1
L26	1
P27	1
V31	1
L32	1
A33	1
Q34	1
E35	1
L36	1
L37	1
L38	1
E41	1
T42	1
E43	1
L44	1
T45	1
E46	1
L47	1
K50	1
F51	1
G52	1
T53	1
G54	1
F57	1
V58	1
D62	1
D63	1
V66	1
R69	1
R70	1
E71	1
Q72	1
D75	1



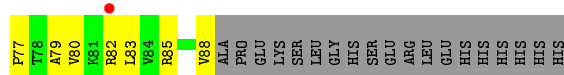
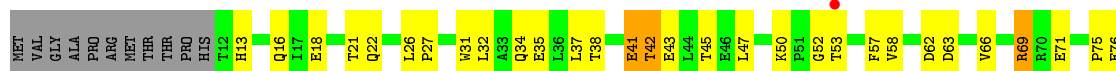
• Molecule 1: SAV_2001



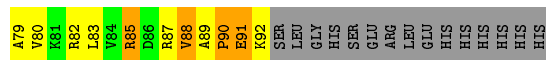
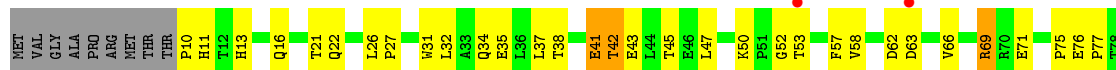
• Molecule 1: SAV_2001



• Molecule 1: SAV_2001



• Molecule 1: SAV_2001



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	48.13Å 158.32Å 48.09Å 90.00° 90.62° 90.00°	Depositor
Resolution (Å)	19.99 – 2.70 28.70 – 2.63	Depositor EDS
% Data completeness (in resolution range)	85.2 (19.99-2.70) 94.1 (28.70-2.63)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.68 (at 2.64Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.230 , 0.264 0.248 , 0.279	Depositor DCC
R_{free} test set	3674 reflections (9.21%)	wwPDB-VP
Wilson B-factor (Å ²)	32.9	Xtriage
Anisotropy	0.423	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 22.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.165 for -l,k,h 0.165 for -h,-k,l 0.458 for l,-k,h	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	5209	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.86% 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	0.48	0/663	0.77	2/903 (0.2%)
1	B	0.46	0/671	0.97	3/913 (0.3%)
1	C	0.49	0/647	1.69	7/881 (0.8%)
1	D	0.48	0/678	0.76	2/923 (0.2%)
1	E	0.48	0/663	0.77	3/903 (0.3%)
1	F	0.45	0/663	0.76	2/903 (0.2%)
1	G	0.46	0/647	0.76	2/881 (0.2%)
1	H	0.48	0/697	0.99	3/949 (0.3%)
All	All	0.47	0/5329	0.98	24/7256 (0.3%)

There are no bond length outliers.

The worst 5 of 24 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	82	ARG	NE-CZ-NH2	-28.28	106.16	120.30
1	C	82	ARG	NE-CZ-NH1	26.14	133.37	120.30
1	H	85	ARG	NE-CZ-NH1	-15.76	112.42	120.30
1	B	85	ARG	NE-CZ-NH1	-15.32	112.64	120.30
1	C	85	ARG	NE-CZ-NH1	-15.32	112.64	120.30

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	648	0	641	24	0
1	B	656	0	653	29	0
1	C	633	0	629	26	0
1	D	663	0	660	28	0
1	E	648	0	641	29	0
1	F	648	0	641	32	0
1	G	633	0	629	23	0
1	H	680	0	675	30	0
All	All	5209	0	5169	211	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 20.

The worst 5 of 211 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:F:34:GLN:HE22	1:H:34:GLN:HE22	0.95	0.94
1:B:34:GLN:HE22	1:D:34:GLN:HE22	0.95	0.91
1:F:34:GLN:HE22	1:H:34:GLN:NE2	1.73	0.87
1:B:34:GLN:HE22	1:D:34:GLN:NE2	1.72	0.86
1:E:34:GLN:HE22	1:G:34:GLN:HE22	1.24	0.84

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	77/107 (72%)	69 (90%)	7 (9%)	1 (1%)	12	30
1	B	78/107 (73%)	68 (87%)	7 (9%)	3 (4%)	3	7
1	C	75/107 (70%)	68 (91%)	6 (8%)	1 (1%)	12	30
1	D	79/107 (74%)	69 (87%)	7 (9%)	3 (4%)	3	7

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	E	77/107 (72%)	67 (87%)	7 (9%)	3 (4%)	3	6
1	F	77/107 (72%)	68 (88%)	8 (10%)	1 (1%)	12	30
1	G	75/107 (70%)	68 (91%)	6 (8%)	1 (1%)	12	30
1	H	81/107 (76%)	71 (88%)	7 (9%)	3 (4%)	3	7
All	All	619/856 (72%)	548 (88%)	55 (9%)	16 (3%)	5	13

5 of 16 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	53	THR
1	B	53	THR
1	C	53	THR
1	D	53	THR
1	E	53	THR

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	69/94 (73%)	61 (88%)	8 (12%)	5	12
1	B	70/94 (74%)	63 (90%)	7 (10%)	7	18
1	C	68/94 (72%)	61 (90%)	7 (10%)	7	16
1	D	71/94 (76%)	60 (84%)	11 (16%)	2	7
1	E	69/94 (73%)	61 (88%)	8 (12%)	5	12
1	F	69/94 (73%)	62 (90%)	7 (10%)	7	17
1	G	68/94 (72%)	61 (90%)	7 (10%)	7	16
1	H	73/94 (78%)	64 (88%)	9 (12%)	4	11
All	All	557/752 (74%)	493 (88%)	64 (12%)	5	13

5 of 64 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	71	GLU
1	E	41	GLU
1	H	69	ARG
1	D	72	GLN
1	D	92	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 23 such sidechains are listed below:

Mol	Chain	Res	Type
1	D	22	GLN
1	E	16	GLN
1	H	16	GLN
1	D	34	GLN
1	E	22	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](#)

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	79/107 (73%)	0.26	4 (5%) 28 26	9, 31, 54, 68	0
1	B	80/107 (74%)	0.24	2 (2%) 57 59	13, 30, 54, 59	0
1	C	77/107 (71%)	0.17	1 (1%) 77 78	13, 32, 53, 56	0
1	D	81/107 (75%)	0.21	3 (3%) 41 41	14, 30, 53, 57	0
1	E	79/107 (73%)	0.25	3 (3%) 40 39	8, 31, 52, 64	0
1	F	79/107 (73%)	0.19	3 (3%) 40 39	15, 31, 53, 58	0
1	G	77/107 (71%)	0.18	2 (2%) 56 57	13, 31, 54, 57	0
1	H	83/107 (77%)	0.22	2 (2%) 59 60	13, 30, 56, 64	0
All	All	635/856 (74%)	0.22	20 (3%) 49 49	8, 31, 54, 68	0

The worst 5 of 20 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	H	53	THR	4.5
1	B	53	THR	4.2
1	D	53	THR	3.9
1	E	53	THR	3.7
1	C	53	THR	3.4

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.