



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 2, 2023 – 01:06 PM EDT

PDB ID : 3Q91
Title : Crystal Structure of Human Uridine Diphosphate Glucose Pyrophosphatase (NUDT14)
Authors : Tresaugues, L.; Siponen, M.I.; Arrowsmith, C.H.; Berglund, H.; Bountra, C.; Collins, R.; Edwards, A.M.; Ekblad, T.; Flodin, S.; Flores, A.; Graslund, S.; Hammarstrom, M.; Johansson, I.; Karlberg, T.; Kol, S.; Kotenyova, T.; Kouznetsova, E.; Moche, M.; Nyman, T.; Persson, C.; Schuler, H.; Schutz, P.; Thorsell, A.G.; Van Der Berg, S.; Wahlberg, E.; Weigelt, J.; Welin, M.; Nordlund, P.; Structural Genomics Consortium (SGC)
Deposited on : 2011-01-07
Resolution : 2.70 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
Xtrriage (Phenix) : 1.13
EDS : 2.35
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)

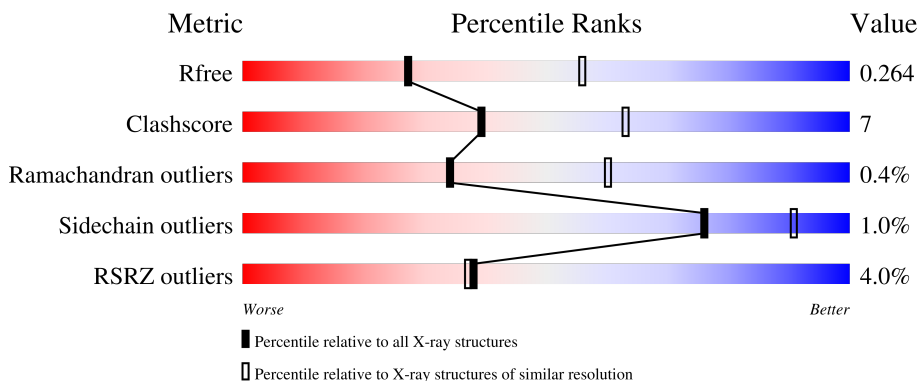
1 Overall quality at a glance

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)

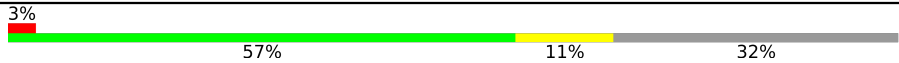
The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. 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	218	
1	B	218	
1	C	218	

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Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
 Validation Pipeline (wwPDB-VP) : 2.35

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Mol	Chain	Length	Quality of chain
1	D	218	 <p>A horizontal bar chart representing the quality of chain. The bar is divided into four segments: a small red segment (3%), a large green segment (57%), a yellow segment (11%), and a grey segment (29%). The percentages are labeled below the bar.</p>

2 Entry composition i

There are 2 unique types of molecules in this entry. The entry contains 4404 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 Uridine diphosphate glucose pyrophosphatase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	145	1116	709	186	217	4	0	0	0
1	C	146	1127	716	190	217	4	0	0	0
1	B	131	1009	642	169	194	4	0	0	0
1	D	149	1147	730	196	217	4	0	0	0

There are 92 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	5	MET	-	expression tag	UNP O95848
A	6	HIS	-	expression tag	UNP O95848
A	7	HIS	-	expression tag	UNP O95848
A	8	HIS	-	expression tag	UNP O95848
A	9	HIS	-	expression tag	UNP O95848
A	10	HIS	-	expression tag	UNP O95848
A	11	HIS	-	expression tag	UNP O95848
A	12	SER	-	expression tag	UNP O95848
A	13	SER	-	expression tag	UNP O95848
A	14	GLY	-	expression tag	UNP O95848
A	15	VAL	-	expression tag	UNP O95848
A	16	ASP	-	expression tag	UNP O95848
A	17	LEU	-	expression tag	UNP O95848
A	18	GLY	-	expression tag	UNP O95848
A	19	THR	-	expression tag	UNP O95848
A	20	GLU	-	expression tag	UNP O95848
A	21	ASN	-	expression tag	UNP O95848
A	22	LEU	-	expression tag	UNP O95848
A	23	TYR	-	expression tag	UNP O95848
A	24	PHE	-	expression tag	UNP O95848
A	25	GLN	-	expression tag	UNP O95848

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Chain	Residue	Modelled	Actual	Comment	Reference
A	26	SER	-	expression tag	UNP O95848
A	27	MET	-	expression tag	UNP O95848
C	5	MET	-	expression tag	UNP O95848
C	6	HIS	-	expression tag	UNP O95848
C	7	HIS	-	expression tag	UNP O95848
C	8	HIS	-	expression tag	UNP O95848
C	9	HIS	-	expression tag	UNP O95848
C	10	HIS	-	expression tag	UNP O95848
C	11	HIS	-	expression tag	UNP O95848
C	12	SER	-	expression tag	UNP O95848
C	13	SER	-	expression tag	UNP O95848
C	14	GLY	-	expression tag	UNP O95848
C	15	VAL	-	expression tag	UNP O95848
C	16	ASP	-	expression tag	UNP O95848
C	17	LEU	-	expression tag	UNP O95848
C	18	GLY	-	expression tag	UNP O95848
C	19	THR	-	expression tag	UNP O95848
C	20	GLU	-	expression tag	UNP O95848
C	21	ASN	-	expression tag	UNP O95848
C	22	LEU	-	expression tag	UNP O95848
C	23	TYR	-	expression tag	UNP O95848
C	24	PHE	-	expression tag	UNP O95848
C	25	GLN	-	expression tag	UNP O95848
C	26	SER	-	expression tag	UNP O95848
C	27	MET	-	expression tag	UNP O95848
B	5	MET	-	expression tag	UNP O95848
B	6	HIS	-	expression tag	UNP O95848
B	7	HIS	-	expression tag	UNP O95848
B	8	HIS	-	expression tag	UNP O95848
B	9	HIS	-	expression tag	UNP O95848
B	10	HIS	-	expression tag	UNP O95848
B	11	HIS	-	expression tag	UNP O95848
B	12	SER	-	expression tag	UNP O95848
B	13	SER	-	expression tag	UNP O95848
B	14	GLY	-	expression tag	UNP O95848
B	15	VAL	-	expression tag	UNP O95848
B	16	ASP	-	expression tag	UNP O95848
B	17	LEU	-	expression tag	UNP O95848
B	18	GLY	-	expression tag	UNP O95848
B	19	THR	-	expression tag	UNP O95848
B	20	GLU	-	expression tag	UNP O95848
B	21	ASN	-	expression tag	UNP O95848

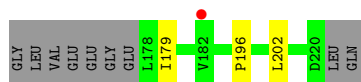
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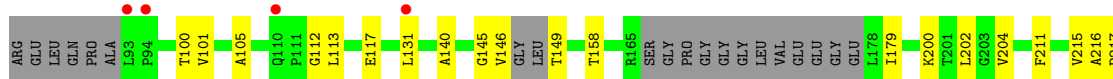
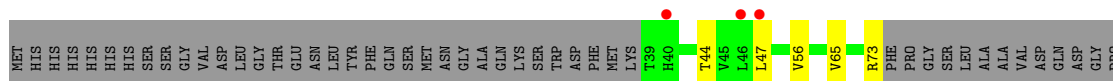
Chain	Residue	Modelled	Actual	Comment	Reference
B	22	LEU	-	expression tag	UNP O95848
B	23	TYR	-	expression tag	UNP O95848
B	24	PHE	-	expression tag	UNP O95848
B	25	GLN	-	expression tag	UNP O95848
B	26	SER	-	expression tag	UNP O95848
B	27	MET	-	expression tag	UNP O95848
D	5	MET	-	expression tag	UNP O95848
D	6	HIS	-	expression tag	UNP O95848
D	7	HIS	-	expression tag	UNP O95848
D	8	HIS	-	expression tag	UNP O95848
D	9	HIS	-	expression tag	UNP O95848
D	10	HIS	-	expression tag	UNP O95848
D	11	HIS	-	expression tag	UNP O95848
D	12	SER	-	expression tag	UNP O95848
D	13	SER	-	expression tag	UNP O95848
D	14	GLY	-	expression tag	UNP O95848
D	15	VAL	-	expression tag	UNP O95848
D	16	ASP	-	expression tag	UNP O95848
D	17	LEU	-	expression tag	UNP O95848
D	18	GLY	-	expression tag	UNP O95848
D	19	THR	-	expression tag	UNP O95848
D	20	GLU	-	expression tag	UNP O95848
D	21	ASN	-	expression tag	UNP O95848
D	22	LEU	-	expression tag	UNP O95848
D	23	TYR	-	expression tag	UNP O95848
D	24	PHE	-	expression tag	UNP O95848
D	25	GLN	-	expression tag	UNP O95848
D	26	SER	-	expression tag	UNP O95848
D	27	MET	-	expression tag	UNP O95848

- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total O 1 1	0	0
2	C	1	Total O 1 1	0	0
2	B	2	Total O 2 2	0	0
2	D	1	Total O 1 1	0	0



- Molecule 1: Uridine diphosphate glucose pyrophosphatase



4 Data and refinement statistics i

Property	Value	Source
Space group	P 2 21 21	Depositor
Cell constants a, b, c, α , β , γ	84.04Å 106.24Å 106.35Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	53.17 – 2.70 53.17 – 2.70	Depositor EDS
% Data completeness (in resolution range)	99.5 (53.17-2.70) 99.5 (53.17-2.70)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.90 (at 2.69Å)	Xtrriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.228 , 0.271 0.225 , 0.264	Depositor DCC
R_{free} test set	1375 reflections (5.15%)	wwPDB-VP
Wilson B-factor (Å ²)	73.7	Xtrriage
Anisotropy	0.457	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 56.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	0.025 for -h,l,k	Xtrriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4404	wwPDB-VP
Average B, all atoms (Å ²)	81.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.10% 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	0.44	0/1138	0.55	0/1548
1	B	0.42	0/1028	0.57	0/1398
1	C	0.45	0/1149	0.55	0/1562
1	D	0.44	0/1170	0.58	0/1592
All	All	0.44	0/4485	0.56	0/6100

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	1116	0	1082	18	0
1	B	1009	0	987	18	0
1	C	1127	0	1097	18	0
1	D	1147	0	1125	16	0
2	A	1	0	0	0	0
2	B	2	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
All	All	4404	0	4291	64	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 7.

All (64) 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:C:58:VAL:HG13	1:C:179:ILE:HG23	1.65	0.77
1:D:65:VAL:HG21	1:D:100:THR:HG21	1.66	0.76
1:B:58:VAL:HG11	1:B:179:ILE:HD12	1.73	0.70
1:B:139:VAL:HG21	1:B:155:MET:CE	2.23	0.69
1:A:70:VAL:HG12	1:A:70:VAL:O	1.93	0.69
1:C:65:VAL:HG21	1:C:100:THR:CG2	2.23	0.67
1:B:60:GLN:HG2	1:B:179:ILE:HG22	1.79	0.64
1:A:59:LYS:HG3	1:A:182:VAL:HG21	1.82	0.62
1:D:101:VAL:HG12	1:D:204:VAL:HG21	1.81	0.62
1:D:113:LEU:HD12	1:D:117:GLU:HB3	1.81	0.61
1:C:205:ILE:HD13	1:D:140:ALA:HB1	1.83	0.60
1:C:201:THR:OG1	1:C:204:VAL:HG23	2.02	0.60
1:C:145:GLY:O	1:C:146:VAL:HG22	2.01	0.59
1:A:186:LEU:HD11	1:A:211:PHE:CE2	2.37	0.59
1:B:116:GLU:CB	1:B:136:LEU:HD12	2.33	0.59
1:A:204:VAL:O	1:A:208:VAL:HG23	2.05	0.56
1:C:58:VAL:CG1	1:C:179:ILE:HG23	2.33	0.56
1:B:46:LEU:HD11	1:B:55:LEU:HD22	1.88	0.55
1:D:65:VAL:HG21	1:D:100:THR:CG2	2.36	0.55
1:A:202:LEU:HD23	1:B:202:LEU:HB3	1.87	0.55
1:D:47:LEU:HD21	1:D:131:LEU:HD11	1.88	0.55
1:A:136:LEU:HD23	1:A:158:THR:HB	1.89	0.55
1:A:131:LEU:HD13	1:A:136:LEU:HD21	1.89	0.54
1:C:202:LEU:HD23	1:D:202:LEU:HD23	1.89	0.54
1:C:65:VAL:HG21	1:C:100:THR:HG21	1.90	0.54
1:C:99:VAL:O	1:C:99:VAL:HG13	2.07	0.53
1:C:47:LEU:HD12	1:C:47:LEU:N	2.24	0.53
1:B:139:VAL:HG21	1:B:155:MET:HE2	1.91	0.52
1:D:179:ILE:O	1:D:179:ILE:HG22	2.10	0.52
1:A:206:PHE:CD2	1:B:202:LEU:HD11	2.46	0.51
1:B:131:LEU:HD22	1:B:158:THR:HG21	1.91	0.51
1:B:116:GLU:HB2	1:B:136:LEU:HD12	1.91	0.51
1:B:139:VAL:HG22	1:B:155:MET:HB3	1.93	0.50
1:A:114:SER:HB3	1:A:116:GLU:HG2	1.95	0.49
1:B:139:VAL:HG21	1:B:155:MET:HE1	1.94	0.49
1:A:141:THR:HG21	1:B:196:PRO:HB3	1.94	0.49
1:D:47:LEU:CD2	1:D:158:THR:HG22	2.43	0.49
1:D:44:THR:HG22	1:D:105:ALA:HB2	1.93	0.48
1:C:65:VAL:HG21	1:C:100:THR:HG23	1.94	0.48
1:D:47:LEU:HB2	1:D:56:VAL:HB	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:49:ASN:HA	1:C:160:VAL:O	2.16	0.46
1:D:211:PHE:CE1	1:D:215:VAL:HG11	2.51	0.46
1:A:215:VAL:O	1:A:219:LEU:HD13	2.16	0.45
1:B:45:VAL:HG13	1:B:104:CYS:HB3	1.99	0.45
1:C:64:ALA:HB2	1:D:149:THR:HA	1.98	0.45
1:A:178:LEU:HD23	1:A:178:LEU:O	2.17	0.44
1:C:58:VAL:HG13	1:C:179:ILE:CG2	2.42	0.44
1:B:108:VAL:HG21	1:B:115:LEU:HD23	2.00	0.44
1:A:206:PHE:CE1	1:A:210:TRP:HB2	2.53	0.44
1:A:70:VAL:O	1:A:70:VAL:CG1	2.64	0.44
1:B:60:GLN:CG	1:B:179:ILE:HG22	2.47	0.43
1:A:47:LEU:HD12	1:A:47:LEU:N	2.33	0.43
1:D:216:ALA:N	1:D:217:PRO:HD2	2.34	0.43
1:A:126:GLU:HG2	1:A:179:ILE:HD13	2.01	0.43
1:B:42:SER:HB3	1:B:107:LEU:HD23	2.01	0.43
1:C:45:VAL:HG22	1:C:47:LEU:HD12	2.01	0.42
1:A:104:CYS:HB2	1:A:127:CYS:SG	2.60	0.42
1:C:99:VAL:O	1:C:99:VAL:CG1	2.68	0.42
1:C:198:ILE:O	1:C:200:LYS:HG3	2.19	0.41
1:A:65:VAL:HG21	1:A:100:THR:CG2	2.50	0.41
1:D:112:GLY:O	1:D:113:LEU:HD22	2.20	0.41
1:D:145:GLY:O	1:D:146:VAL:C	2.60	0.41
1:C:216:ALA:N	1:C:217:PRO:HD2	2.36	0.40
1:B:47:LEU:HD13	1:B:56:VAL:HB	2.03	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	137/218 (63%)	128 (93%)	8 (6%)	1 (1%)	22 46

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	123/218 (56%)	118 (96%)	5 (4%)	0	100	100
1	C	138/218 (63%)	133 (96%)	4 (3%)	1 (1%)	22	46
1	D	141/218 (65%)	134 (95%)	7 (5%)	0	100	100
All	All	539/872 (62%)	513 (95%)	24 (4%)	2 (0%)	34	60

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	219	LEU
1	A	70	VAL

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	121/179 (68%)	121 (100%)	0	100	100
1	B	111/179 (62%)	110 (99%)	1 (1%)	78	92
1	C	122/179 (68%)	120 (98%)	2 (2%)	62	85
1	D	124/179 (69%)	122 (98%)	2 (2%)	62	85
All	All	478/716 (67%)	473 (99%)	5 (1%)	76	91

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	180	GLU
1	C	182	VAL
1	B	139	VAL
1	D	73	ARG
1	D	200	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	60	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	145/218 (66%)	0.46	4 (2%) 53 54	52, 77, 118, 162	0
1	B	131/218 (60%)	0.53	7 (5%) 26 25	52, 77, 115, 131	0
1	C	146/218 (66%)	0.51	5 (3%) 45 45	47, 78, 120, 151	0
1	D	149/218 (68%)	0.46	7 (4%) 31 30	57, 80, 122, 197	0
All	All	571/872 (65%)	0.49	23 (4%) 38 37	47, 78, 120, 197	0

All (23) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	70	VAL	6.2
1	A	70	VAL	4.7
1	D	93	LEU	3.3
1	D	47	LEU	3.2
1	B	113	LEU	2.8
1	A	178	LEU	2.7
1	D	40	HIS	2.5
1	B	146	VAL	2.5
1	B	111	PRO	2.5
1	C	150	GLY	2.4
1	B	112	GLY	2.3
1	D	131	LEU	2.3
1	A	131	LEU	2.3
1	C	143	TRP	2.2
1	B	40	HIS	2.2
1	B	131	LEU	2.2
1	D	94	PRO	2.2
1	A	186	LEU	2.2
1	D	46	LEU	2.2
1	D	110	GLN	2.2
1	C	115	LEU	2.1

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Mol	Chain	Res	Type	RSRZ
1	C	131	LEU	2.1
1	B	182	VAL	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 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.