



wwPDB X-ray Structure Validation Summary Report ⓘ

Jan 31, 2016 – 10:42 PM GMT

PDB ID : 1USV
Title : THE STRUCTURE OF THE COMPLEX BETWEEN AHA1 AND HSP90
Authors : Meyer, P.; Roe, S.M.; Pearl, L.H.
Deposited on : 2003-12-01
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
<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

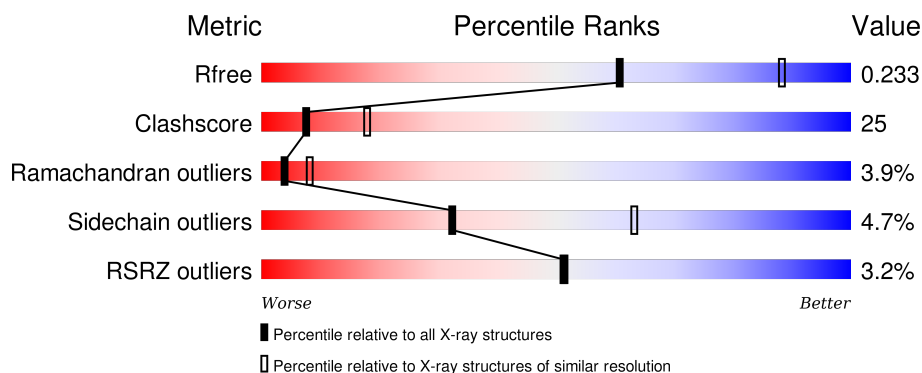
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 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}	91344	2103 (2.70-2.70)
Clashscore	102246	2422 (2.70-2.70)
Ramachandran outliers	100387	2382 (2.70-2.70)
Sidechain outliers	100360	2382 (2.70-2.70)
RSRZ outliers	91569	2107 (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. 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	260	<div> <div>3%</div> <div>56% 36% 5% .</div> </div>
1	C	260	<div> <div>%</div> <div>61% 30% . 5%</div> </div>
1	E	260	<div> <div>4%</div> <div>53% 39% . .</div> </div>
1	G	260	<div> <div>%</div> <div>50% 41% 5% 5%</div> </div>
2	B	170	<div> <div>4%</div> <div>42% 38% . 16%</div> </div>

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Mol	Chain	Length	Quality of chain
2	D	170	<div><div></div><div>3%</div><div>49%</div><div>30%</div><div>5%</div><div>16%</div></div>
2	F	170	<div><div></div><div>6%</div><div>39%</div><div>34%</div><div>12%</div><div>•</div><div>14%</div></div>
2	H	170	<div><div></div><div>4%</div><div>48%</div><div>34%</div><div>••</div><div>17%</div></div>

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 12783 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 HEAT SHOCK PROTEIN HSP82.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	251	Total	C	N	O	S	0	0	0
			2080	1345	338	394	3			
1	C	247	Total	C	N	O	S	0	0	0
			2041	1320	332	387	2			
1	E	251	Total	C	N	O	S	0	0	0
			2078	1342	339	394	3			
1	G	248	Total	C	N	O	S	0	0	0
			2050	1325	335	387	3			

- Molecule 2 is a protein called AHA1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	143	Total	C	N	O	S	0	0	0
			1124	719	185	218	2			
2	D	142	Total	C	N	O	S	0	0	0
			1107	707	183	215	2			
2	F	146	Total	C	N	O	S	0	0	1
			1146	735	191	218	2			
2	H	141	Total	C	N	O	S	0	0	0
			1099	703	181	213	2			

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	16	Total	O	0	0
			16	16		
3	B	4	Total	O	0	0
			4	4		
3	C	13	Total	O	0	0
			13	13		
3	D	2	Total	O	0	0
			2	2		

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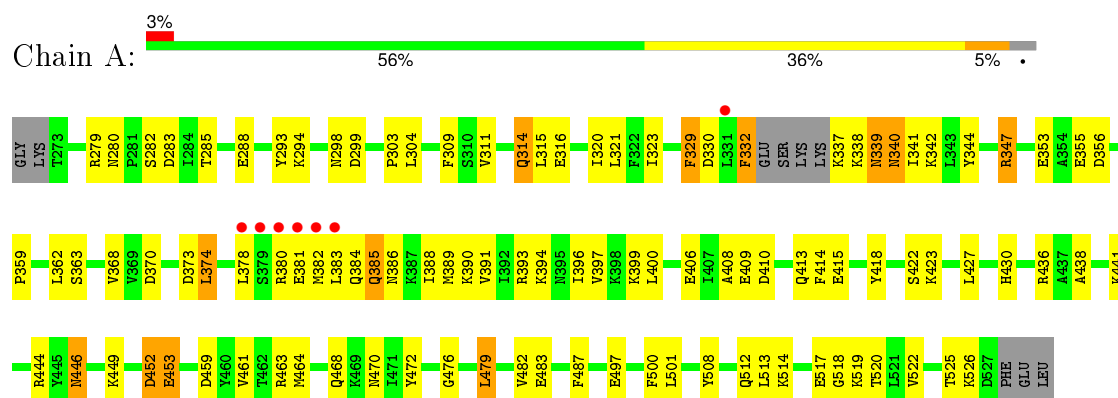
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	E	7	Total 7	O 7	0	0
3	F	2	Total 2	O 2	0	0
3	G	11	Total 11	O 11	0	0
3	H	3	Total 3	O 3	0	0

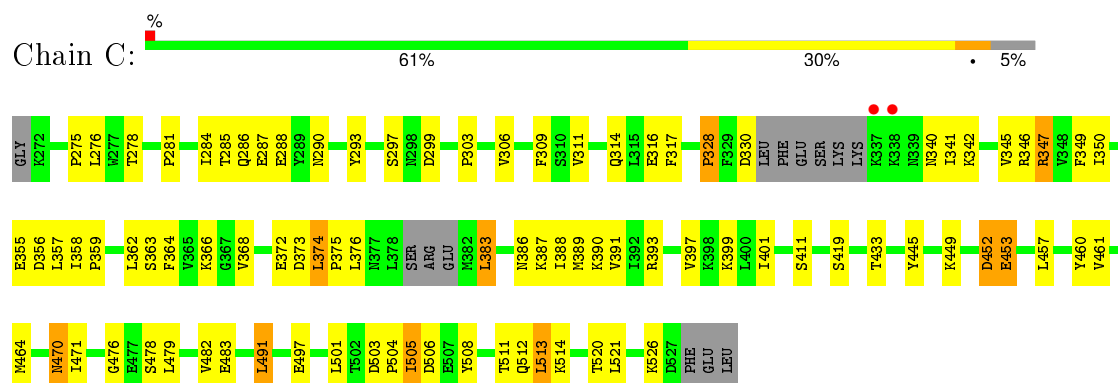
3 Residue-property plots

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 ($RSRZ > 2$). 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.

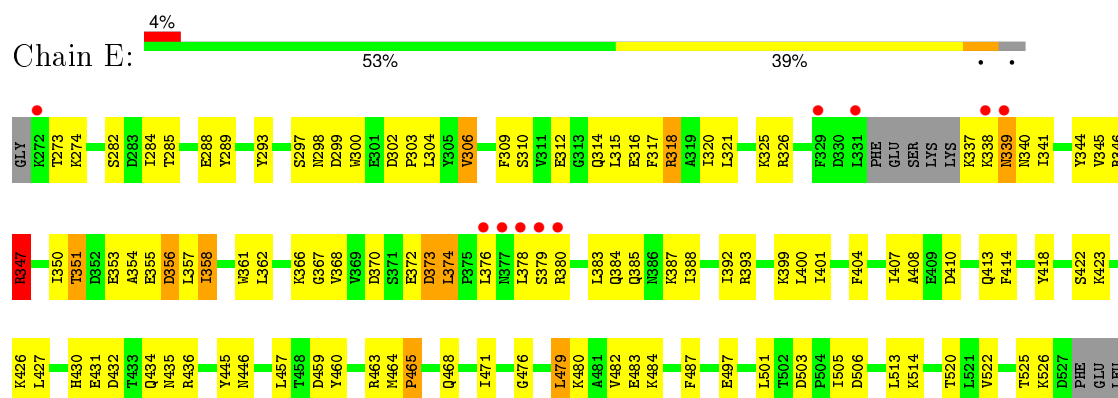
• Molecule 1: HEAT SHOCK PROTEIN HSP82



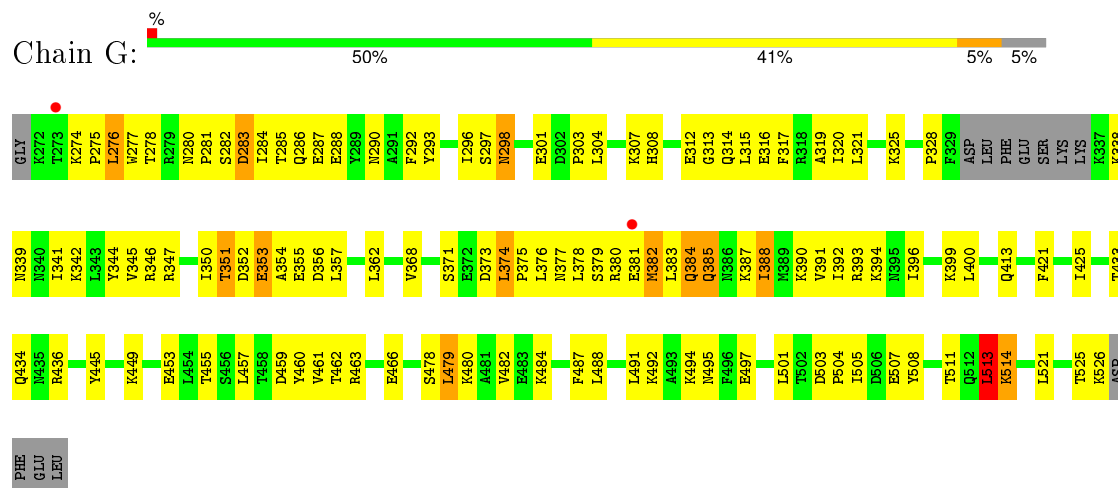
• Molecule 1: HEAT SHOCK PROTEIN HSP82



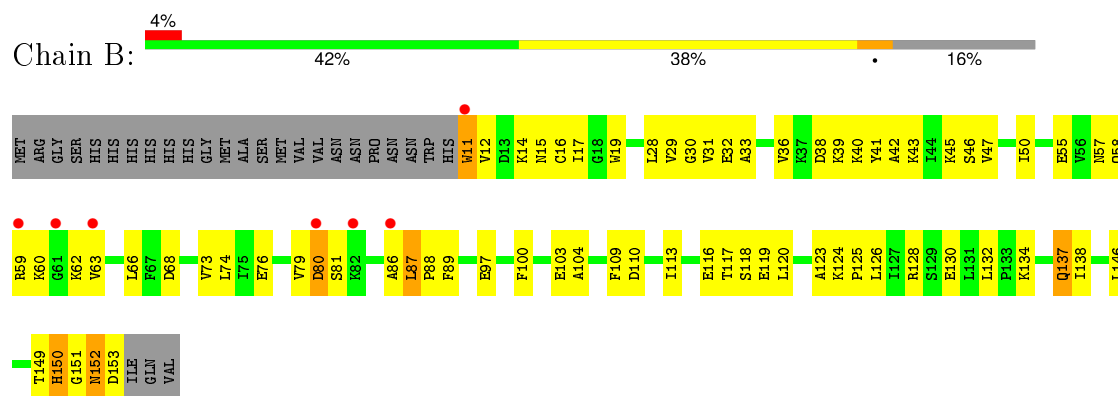
• Molecule 1: HEAT SHOCK PROTEIN HSP82

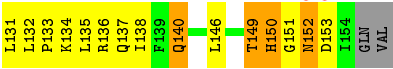


- Molecule 1: HEAT SHOCK PROTEIN HSP82

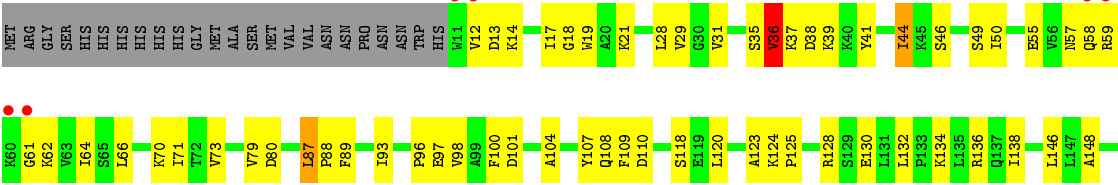


- Molecule 2: AHA1





● Molecule 2: AHA1



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	58.40 Å 207.93 Å 84.45 Å 90.00° 97.68° 90.00°	Depositor
Resolution (Å)	34.26 – 2.70 34.26 – 2.70	Depositor EDS
% Data completeness (in resolution range)	97.8 (34.26-2.70) 97.9 (34.26-2.70)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.85 (at 2.68 Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.229 , 0.298 0.231 , 0.233	Depositor DCC
R_{free} test set	2720 reflections (5.37%)	DCC
Wilson B-factor (Å ²)	51.6	Xtriage
Anisotropy	0.444	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 45.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Outliers	0 of 53401 reflections	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	12783	wwPDB-VP
Average B, all atoms (Å ²)	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.81% 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, 0.375 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	0.46	0/2124	0.64	0/2867
1	C	0.48	0/2083	0.66	0/2811
1	E	0.44	0/2121	0.65	0/2862
1	G	0.44	0/2093	0.64	0/2825
2	B	0.40	0/1144	0.65	0/1537
2	D	0.43	0/1125	0.67	0/1510
2	F	0.39	0/1169	0.60	0/1573
2	H	0.41	0/1117	0.67	0/1499
All	All	0.44	0/12976	0.65	0/17484

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	2080	0	2083	91	0
1	C	2041	0	2044	92	0
1	E	2078	0	2087	112	0
1	G	2050	0	2057	116	0
2	B	1124	0	1132	74	0
2	D	1107	0	1120	51	0
2	F	1146	0	1147	79	0
2	H	1099	0	1114	51	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	16	0	0	0	0
3	B	4	0	0	0	0
3	C	13	0	0	1	0
3	D	2	0	0	0	0
3	E	7	0	0	1	0
3	F	2	0	0	0	0
3	G	11	0	0	0	0
3	H	3	0	0	0	0
All	All	12783	0	12784	647	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 25.

The worst 5 of 647 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:G:280:ASN:HD22	1:G:283:ASP:HB2	1.13	1.11
1:G:380:ARG:HG2	1:G:381:GLU:H	1.23	1.03
1:C:390:LYS:HA	1:C:393:ARG:HH12	1.25	0.98
1:C:386:ASN:HD22	1:C:389:MET:H	1.05	0.97
2:F:78:HIS:HB2	2:F:86:ALA:HB1	1.47	0.95

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	247/260 (95%)	219 (89%)	21 (8%)	7 (3%)	6	15
1	C	241/260 (93%)	214 (89%)	19 (8%)	8 (3%)	5	11
1	E	247/260 (95%)	207 (84%)	33 (13%)	7 (3%)	6	15

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	G	244/260 (94%)	200 (82%)	34 (14%)	10 (4%)	3	7
2	B	141/170 (83%)	111 (79%)	27 (19%)	3 (2%)	9	23
2	D	140/170 (82%)	121 (86%)	15 (11%)	4 (3%)	6	14
2	F	144/170 (85%)	102 (71%)	25 (17%)	17 (12%)	0	0
2	H	139/170 (82%)	117 (84%)	18 (13%)	4 (3%)	6	14
All	All	1543/1720 (90%)	1291 (84%)	192 (12%)	60 (4%)	4	8

5 of 60 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	385	GLN
1	A	453	GLU
1	C	341	ILE
2	D	36	VAL
2	D	121	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	232/240 (97%)	222 (96%)	10 (4%)	35	66
1	C	227/240 (95%)	216 (95%)	11 (5%)	31	62
1	E	232/240 (97%)	222 (96%)	10 (4%)	35	66
1	G	228/240 (95%)	216 (95%)	12 (5%)	28	57
2	B	125/149 (84%)	120 (96%)	5 (4%)	38	69
2	D	123/149 (83%)	117 (95%)	6 (5%)	31	61
2	F	126/149 (85%)	118 (94%)	8 (6%)	22	48
2	H	122/149 (82%)	117 (96%)	5 (4%)	37	69
All	All	1415/1556 (91%)	1348 (95%)	67 (5%)	32	63

5 of 67 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	D	140	GLN
1	E	356	ASP
1	G	521	LEU
2	D	152	ASN
1	E	318	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 51 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	446	ASN
1	E	280	ASN
1	G	405	ASN
2	D	108	GLN
1	E	339	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 ⓘ

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	251/260 (96%)	-0.21	7 (2%) 56 57	22, 44, 81, 114	0
1	C	247/260 (95%)	-0.28	2 (0%) 87 88	19, 40, 67, 87	0
1	E	251/260 (96%)	-0.15	10 (3%) 42 41	25, 45, 89, 120	0
1	G	248/260 (95%)	-0.11	2 (0%) 87 88	22, 51, 97, 115	0
2	B	143/170 (84%)	-0.07	7 (4%) 33 32	25, 56, 95, 103	0
2	D	142/170 (83%)	-0.07	5 (3%) 48 48	30, 56, 95, 101	0
2	F	146/170 (85%)	0.31	11 (7%) 17 15	38, 80, 109, 114	0
2	H	141/170 (82%)	-0.12	6 (4%) 39 38	41, 59, 92, 109	0
All	All	1569/1720 (91%)	-0.11	50 (3%) 51 51	19, 50, 98, 120	0

The worst 5 of 50 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	H	61	GLY	5.4
2	B	61	GLY	4.6
1	A	331	LEU	4.6
2	H	11	TRP	4.3
1	A	382	MET	4.3

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.