



Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 12:37 AM GMT

PDB ID : 2QA7
Title : Crystal structure of Huntingtin-interactingprotein 1 (HIP1) coiled-coil domain
with a basic surface suitable for HIP-protein interactor (HIPPI)
Authors : Niu, Q.; Ybe, J.A.
Deposited on : 2007-06-14
Resolution : 2.80 Å(reported)

This is a full wwPDB validation 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/ValidationPDFNotes.html>

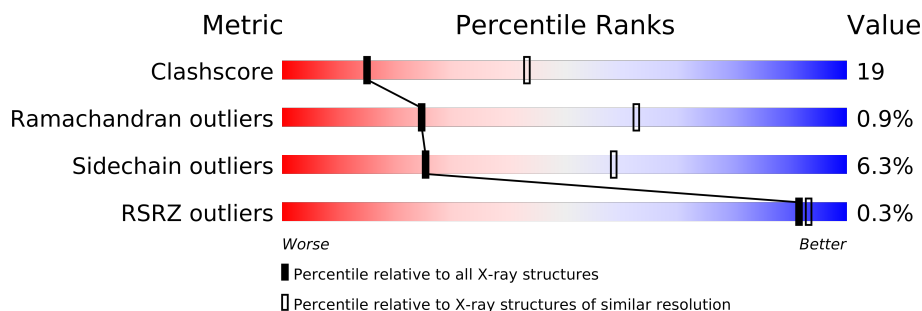
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 2.80 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))
Clashscore	79885	2295 (2.80-2.80)
Ramachandran outliers	78287	2252 (2.80-2.80)
Sidechain outliers	78261	2254 (2.80-2.80)
RSRZ outliers	66119	1802 (2.80-2.80)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	114	
1	B	114	
1	C	114	
1	D	114	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 2814 atoms, of which 0 are hydrogen and 0 are deuterium.

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 Huntingtin-interactingprotein 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	99	Total	C	N	O	S	0	0	0
			821	494	157	168	2			
1	B	70	Total	C	N	O	S	0	0	0
			572	350	103	117	2			
1	C	102	Total	C	N	O	S	0	0	0
			847	512	162	171	2			
1	D	70	Total	C	N	O	S	0	0	0
			572	350	103	117	2			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	368	GLY	-	EXPRESSION TAG	UNP O00291
A	369	SER	-	EXPRESSION TAG	UNP O00291
B	368	GLY	-	EXPRESSION TAG	UNP O00291
B	369	SER	-	EXPRESSION TAG	UNP O00291
C	368	GLY	-	EXPRESSION TAG	UNP O00291
C	369	SER	-	EXPRESSION TAG	UNP O00291
D	368	GLY	-	EXPRESSION TAG	UNP O00291
D	369	SER	-	EXPRESSION TAG	UNP O00291

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	D	2	Total	O	0	0
			2	2		

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.

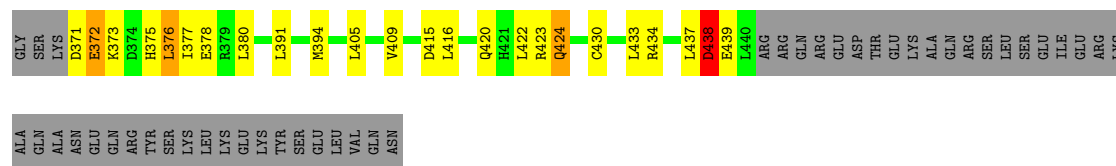
• Molecule 1: Huntingtin-interactingprotein 1

Chain A: 



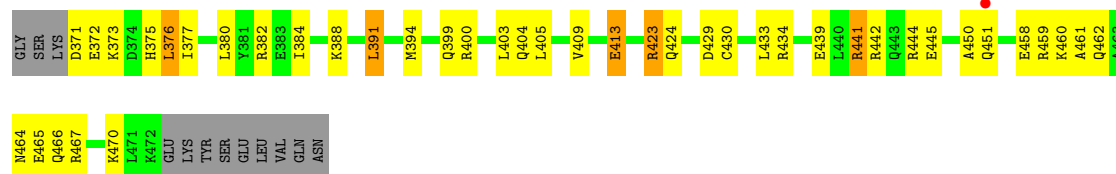
• Molecule 1: Huntingtin-interactingprotein 1

Chain B: 



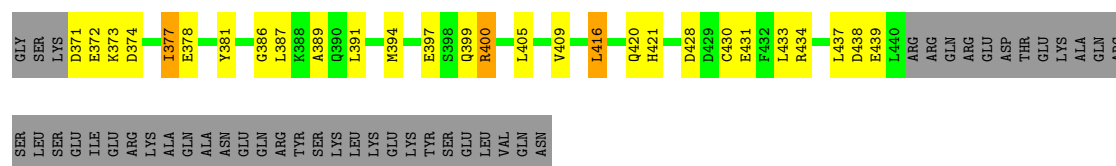
• Molecule 1: Huntingtin-interactingprotein 1

Chain C: 



• Molecule 1: Huntingtin-interactingprotein 1

Chain D: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 43	Depositor
Cell constants a, b, c, α , β , γ	72.90Å 72.90Å 106.50Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.80 43.00 – 2.70	Depositor EDS
% Data completeness (in resolution range)	98.6 (30.00-2.80) 99.2 (43.00-2.70)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.13 (at 2.69Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.265 , 0.324 0.256 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	54.6	Xtriage
Anisotropy	0.523	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 40.6	EDS
Estimated twinning fraction	0.478 for h,-k,-l	Xtriage
L-test for twinning	$\langle L \rangle = 0.54$, $\langle L^2 \rangle = 0.38$	Xtriage
Outliers	0 of 15317 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	2814	wwPDB-VP
Average B, all atoms (Å ²)	62.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.66% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

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	0.58	2/826 (0.2%)	0.63	1/1102 (0.1%)
1	B	0.40	0/576	0.55	0/771
1	C	0.46	1/852 (0.1%)	0.59	0/1135
1	D	0.40	0/576	0.57	0/771
All	All	0.48	3/2830 (0.1%)	0.59	1/3779 (0.0%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	413	GLU	CG-CD	8.94	1.65	1.51
1	A	413	GLU	CB-CG	7.11	1.65	1.52
1	C	413	GLU	CG-CD	6.16	1.61	1.51

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	413	GLU	OE1-CD-OE2	-7.30	114.55	123.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Close contacts

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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	821	0	803	38	1

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	572	0	559	38	0
1	C	847	0	840	41	1
1	D	572	0	559	19	0
2	D	2	0	0	1	0
All	All	2814	0	2761	106	2

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 19.

All (106) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:371:ASP:HB3	1:B:375:HIS:CE1	2.01	0.95
1:A:443:GLN:HA	1:C:382:ARG:HH12	1.33	0.91
1:A:443:GLN:HA	1:C:382:ARG:NH1	1.92	0.83
1:A:443:GLN:HA	1:C:382:ARG:HH22	1.44	0.82
1:B:372:GLU:HG3	1:B:373:LYS:H	1.42	0.81
1:C:399:GLN:O	1:C:403:LEU:HD23	1.79	0.81
1:C:400:ARG:CZ	1:C:404:GLN:HE22	1.93	0.81
1:B:372:GLU:HG3	1:B:373:LYS:N	1.98	0.77
1:A:443:GLN:HA	1:C:382:ARG:NH2	1.99	0.77
1:D:416:LEU:O	1:D:420:GLN:HB2	1.85	0.76
1:C:391:LEU:HD23	1:D:391:LEU:HD13	1.67	0.75
1:A:371:ASP:HA	1:A:374:ASP:OD2	1.86	0.75
1:C:461:ALA:O	1:C:464:ASN:HB3	1.87	0.74
1:C:371:ASP:OD2	1:C:375:HIS:NE2	2.19	0.73
1:A:437:LEU:HD23	1:B:437:LEU:HD13	1.73	0.71
1:A:430:CYS:HB3	1:B:433:LEU:HD12	1.71	0.71
1:B:372:GLU:CG	1:B:373:LYS:H	2.01	0.70
1:B:430:CYS:HB3	1:B:434:ARG:HH21	1.57	0.69
1:D:397:GLU:OE2	1:D:400:ARG:HD3	1.92	0.69
1:A:443:GLN:HA	1:C:382:ARG:CZ	2.22	0.69
1:C:400:ARG:HG2	1:C:404:GLN:HE21	1.58	0.69
1:A:443:GLN:CA	1:C:382:ARG:HH22	2.06	0.69
1:B:371:ASP:OD1	1:B:375:HIS:NE2	2.24	0.68
1:A:443:GLN:CA	1:C:382:ARG:HH12	2.06	0.68
1:A:395:LYS:HG2	1:B:394:MET:HE2	1.77	0.66
1:B:416:LEU:O	1:B:420:GLN:HB2	1.96	0.66
1:C:400:ARG:NH1	1:C:404:GLN:HE22	1.93	0.65
1:D:428:ASP:O	1:D:431:GLU:HB3	1.97	0.65
1:A:436:GLU:O	1:A:440:LEU:HD23	1.96	0.65
1:C:371:ASP:CG	1:C:375:HIS:HE2	2.01	0.64
1:D:421:HIS:CE1	2:D:1:HOH:O	2.50	0.64

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:395:LYS:HG2	1:B:394:MET:CE	2.28	0.63
1:D:405:LEU:O	1:D:409:VAL:HG23	1.99	0.63
1:B:375:HIS:HA	1:B:378:GLU:HG2	1.82	0.61
1:C:391:LEU:HD23	1:D:391:LEU:CD1	2.29	0.61
1:A:395:LYS:O	1:A:399:GLN:HG2	2.03	0.59
1:C:400:ARG:HG2	1:C:404:GLN:NE2	2.19	0.58
1:B:423:ARG:HH22	1:B:424:GLN:HG2	1.68	0.58
1:B:371:ASP:O	1:B:372:GLU:HG2	2.05	0.56
1:C:371:ASP:CG	1:C:375:HIS:NE2	2.59	0.56
1:B:437:LEU:O	1:B:439:GLU:N	2.39	0.55
1:C:465:GLU:C	1:C:467:ARG:H	2.10	0.55
1:A:415:ASP:O	1:A:418:GLU:HB3	2.06	0.55
1:A:451:GLN:O	1:A:455:SER:HB2	2.07	0.55
1:B:423:ARG:NH2	1:B:424:GLN:HG2	2.21	0.54
1:A:419:GLN:OE1	1:B:420:GLN:HA	2.08	0.53
1:A:430:CYS:HB3	1:B:433:LEU:CD1	2.38	0.53
1:B:375:HIS:O	1:B:378:GLU:HG2	2.08	0.53
1:A:400:ARG:HH11	1:A:400:ARG:HG2	1.73	0.53
1:C:459:ARG:C	1:C:461:ALA:H	2.13	0.53
1:D:373:LYS:O	1:D:377:ILE:HG23	2.10	0.52
1:B:371:ASP:CB	1:B:375:HIS:CE1	2.87	0.52
1:C:439:GLU:O	1:C:442:ARG:HB3	2.09	0.52
1:A:391:LEU:HA	1:B:391:LEU:HD21	1.91	0.51
1:D:437:LEU:C	1:D:439:GLU:H	2.15	0.50
1:B:371:ASP:O	1:B:372:GLU:CB	2.58	0.50
1:C:394:MET:HG2	1:D:394:MET:HE3	1.94	0.49
1:A:412:LEU:O	1:A:413:GLU:C	2.51	0.49
1:D:386:GLY:O	1:D:389:ALA:HB3	2.12	0.49
1:A:394:MET:HE2	1:B:394:MET:SD	2.52	0.49
1:B:405:LEU:O	1:B:409:VAL:HG23	2.12	0.49
1:C:384:ILE:O	1:C:388:LYS:HG3	2.12	0.49
1:A:376:LEU:O	1:A:380:LEU:HG	2.13	0.49
1:D:371:ASP:OD2	1:D:372:GLU:N	2.46	0.48
1:C:462:GLN:C	1:C:464:ASN:H	2.16	0.48
1:A:412:LEU:C	1:A:414:ALA:N	2.67	0.48
1:A:437:LEU:HD23	1:B:437:LEU:CD1	2.43	0.48
1:C:373:LYS:O	1:C:377:ILE:HG12	2.13	0.48
1:C:466:GLN:HG2	1:C:466:GLN:O	2.13	0.48
1:A:444:ARG:NE	1:A:444:ARG:HA	2.29	0.47
1:B:371:ASP:O	1:B:372:GLU:HB3	2.15	0.47
1:A:458:GLU:C	1:A:460:LYS:H	2.17	0.47
1:C:429:ASP:O	1:C:433:LEU:HG	2.15	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:457:ILE:HG12	1:D:372:GLU:HG2	1.95	0.47
1:C:394:MET:HB3	1:D:394:MET:CE	2.44	0.47
1:C:441:ARG:O	1:C:445:GLU:HB2	2.14	0.47
1:B:430:CYS:HB3	1:B:434:ARG:NH2	2.27	0.46
1:A:437:LEU:CD2	1:B:437:LEU:HD13	2.45	0.46
1:C:470:LYS:O	1:C:470:LYS:HG2	2.15	0.46
1:C:372:GLU:OE1	1:C:372:GLU:HA	2.16	0.46
1:B:373:LYS:O	1:B:377:ILE:HG12	2.15	0.46
1:A:394:MET:HE1	1:B:394:MET:HB3	1.98	0.46
1:A:395:LYS:CG	1:B:394:MET:CE	2.95	0.45
1:D:377:ILE:CD1	1:D:381:TYR:CE1	2.99	0.45
1:D:371:ASP:HB3	1:D:374:ASP:OD2	2.16	0.45
1:A:447:THR:HG22	1:A:451:GLN:HE21	1.82	0.44
1:C:459:ARG:C	1:C:461:ALA:N	2.72	0.43
1:C:423:ARG:HH12	1:C:424:GLN:HG2	1.83	0.43
1:C:376:LEU:HD22	1:C:380:LEU:HG	2.00	0.43
1:C:434:ARG:NH1	1:D:433:LEU:HD11	2.35	0.42
1:A:423:ARG:HD2	1:B:422:LEU:HB3	2.00	0.42
1:B:437:LEU:O	1:B:438:ASP:C	2.59	0.41
1:A:391:LEU:HD12	1:B:391:LEU:HD23	2.02	0.41
1:C:450:ALA:O	1:C:451:GLN:C	2.59	0.41
1:C:377:ILE:HD11	1:D:377:ILE:HG22	2.02	0.41
1:C:405:LEU:O	1:C:409:VAL:HG23	2.20	0.41
1:A:410:SER:HA	1:A:413:GLU:HB2	2.03	0.41
1:B:376:LEU:O	1:B:380:LEU:HG	2.20	0.41
1:B:371:ASP:C	1:B:372:GLU:HG2	2.41	0.41
1:D:377:ILE:HG13	1:D:378:GLU:N	2.35	0.41
1:C:400:ARG:NH1	1:C:404:GLN:NE2	2.64	0.41
1:A:430:CYS:N	1:B:430:CYS:SG	2.94	0.41
1:A:464:ASN:O	1:A:467:ARG:HB2	2.21	0.40
1:A:390:GLN:NE2	1:B:391:LEU:HD13	2.37	0.40
1:C:444:ARG:HG3	1:C:445:GLU:N	2.37	0.40
1:C:460:LYS:HG3	1:C:460:LYS:O	2.21	0.40

All (2) 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	Distance(Å)	Clash(Å)
1:C:371:ASP:OD2	1:C:413:GLU:OE2[3_654]	1.97	0.23
1:A:375:HIS:NE2	1:A:413:GLU:OE1[4_575]	2.10	0.10

5.3 Torsion angles

5.3.1 Protein backbone ⓘ

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	97/114 (85%)	91 (94%)	6 (6%)	0	100	100
1	B	68/114 (60%)	60 (88%)	6 (9%)	2 (3%)	7	23
1	C	100/114 (88%)	90 (90%)	10 (10%)	0	100	100
1	D	68/114 (60%)	64 (94%)	3 (4%)	1 (2%)	15	46
All	All	333/456 (73%)	305 (92%)	25 (8%)	3 (1%)	25	63

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	372	GLU
1	B	438	ASP
1	D	438	ASP

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	88/102 (86%)	86 (98%)	2 (2%)	63	92
1	B	62/102 (61%)	58 (94%)	4 (6%)	24	57
1	C	91/102 (89%)	85 (93%)	6 (7%)	24	56
1	D	62/102 (61%)	55 (89%)	7 (11%)	9	24
All	All	303/408 (74%)	284 (94%)	19 (6%)	25	59

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	443	GLN
1	A	455	SER
1	B	376	LEU
1	B	415	ASP
1	B	424	GLN
1	B	438	ASP
1	C	376	LEU
1	C	391	LEU
1	C	423	ARG
1	C	430	CYS
1	C	441	ARG
1	C	458	GLU
1	D	377	ILE
1	D	387	LEU
1	D	399	GLN
1	D	400	ARG
1	D	416	LEU
1	D	430	CYS
1	D	434	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	390	GLN
1	A	451	GLN
1	A	462	GLN
1	B	390	GLN
1	B	399	GLN
1	C	404	GLN
1	C	424	GLN
1	D	393	ASN

5.3.3 RNA ⓘ

There are no RNA chains in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

There are no ligands in this entry.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

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	99/114 (86%)	0.01	0 100 100	34, 58, 109, 119	0
1	B	70/114 (61%)	-0.20	0 100 100	30, 58, 76, 119	0
1	C	102/114 (89%)	-0.02	1 (0%) 79 79	37, 58, 106, 134	0
1	D	70/114 (61%)	-0.16	0 100 100	37, 55, 80, 110	0
All	All	341/456 (74%)	-0.08	1 (0%) 91 93	30, 58, 102, 134	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	451	GLN	2.2

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.