

Full wwPDB X-ray Structure Validation Report (i)

Aug 28, 2023 – 02:44 PM EDT

PDB ID : 3LRH

Title: Structure of anti-huntingtin VL domain in complex with huntingtin peptide

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Deposited on : 2010-02-11

Resolution : 2.60 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 (i) 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 (1)) were used in the production of this report:

 $\begin{array}{ccc} \text{MolProbity} & : & 4.02\text{b-}467 \\ \text{Xtriage (Phenix)} & : & 1.13 \end{array}$

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) roteins) : Engh & Huber (2001)

Ideal geometry (proteins) : Engh & Huber (2001) Ideal geometry (DNA, RNA) : Parkinson et al. (1996)

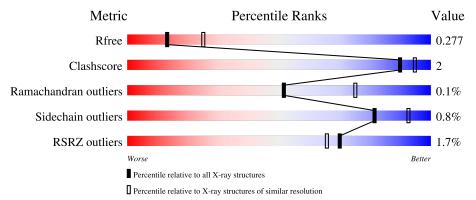
Validation Pipeline (wwPDB-VP) : 2.35

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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 $(\# \mathrm{Entries})$	$\begin{array}{c} {\rm Similar\ resolution} \\ (\#{\rm Entries},{\rm resolution\ range}({\rm \AA})) \end{array}$
R_{free}	130704	3163 (2.60-2.60)
Clashscore	141614	3518 (2.60-2.60)
Ramachandran outliers	138981	3455 (2.60-2.60)
Sidechain outliers	138945	3455 (2.60-2.60)
RSRZ outliers	127900	3104 (2.60-2.60)

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 <=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	121	87%	10%
1	С	121	84% 6%	10%
1	Е	121	88% 5	% 7%
1	G	121	83% 7%	10%
1	I	121	83% 7%	



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Mol	Chain	Length	Quality of chain	
1	K	121	85% 5%	10%
1	M	121	87%	9%
1	О	121	83% 8%	8%
2	В	14	100%	
2	D	14	93%	7%
2	F	14	93%	7%
2	Н	14	14%	
2	J	14	7% 100%	
2	L	14	7% 100%	
2	N	14	93%	7%
2	Р	14	100%	



2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 7652 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 anti-huntingtin VL domain.

Mol	Chain	Residues		At	oms			ZeroOcc	AltConf	Trace
1	A	109	Total	С	N	О	S	0	0	0
1	A	109	807	499	138	169	1	U	U	0
1	С	109	Total	С	N	О	S	0	0	0
1		109	807	499	138	169	1	0	U	0
1	Е	112	Total	С	N	О	S	0	0	0
1	l Li	112	828	510	142	175	1	0	0	
1	G	109	Total	С	N	N O S	0	0	0	
1	G	109	807	499	138	169	1	0	U	U
1	I	110	Total	С	N	О	S	0	0	0
1	1	110	813	502	139	171	1	0	0	
1	K	109	Total	С	N	О	S	0	0	0
1	IX	109	807	499	138	169	1	0	0	0
1	M	110	Total	С	N	О	S	0	0	0
1	IVI	110	813	502	139	171	1	0	0	0
1	O	111	Total	С	N	О	S	0	0	0
1		111	822	507	141	173	1	U		0

• Molecule 2 is a protein called Huntingtin.

Mol	Chain	Residues		Ato	oms			ZeroOcc	AltConf	Trace
2	В	14	Total	С	N	О	S	0	0	0
2	Ъ	14	118	77	18	22	1	0		U
2	D	14	Total	С	N	О	S	0	0	0
2	D	14	118	77	18	22	1	U	U	U
2	F	14	Total	С	N	Ο	S	0	0	0
2	I.	14	118	77	18	22	1	0		U
2	Н	14	Total	С	N	Ο	S	0	0	0
2	11	14	118	77	18	22	1	0		U
2	J	14	Total	С	N	О	S	0	0	0
2	J	14	118	77	18	22	1	0		U
2	L	14	Total	С	N	О	S	0	0	0
	П	14	118	77	18	22	1	U		U



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M	ol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
6)	N	1.4	Total	С	N	О	S	0	0	0
4	_	1N	14	118	77	18	22	1	U	U	U
6)	D	1.4	Total	С	N	О	S	0	0	0
4	4	Г	14	118	77	18	22	1	U	U	U

• Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf	
3	A	21	Total O	0	0	
	11	21	21 21	0	Ŭ .	
3	В	4	Total O	0	0	
5	D	4	4 4		U	
3	\mathbf{C}	16	Total O	0	0	
5)	10	16 16		U	
3	D	2	Total O	0	0	
	D	2	2 2	0	U	
3	E	20	Total O	0	0	
	ם	20	20 20		O	
3	F	1	Total O	0	0	
3	T.	1	1 1	U	U	
3	G	24	Total O	0	0	
	J	24	24 24	0	U	
3	Н	1	Total O	0	0	
	11	1	1 1	U	U	
3	I	22	Total O	0	0	
	1	22	22 22	0	0	
3	J	1	Total O	0	0	
5	5	1	1 1		U	
3	K	35	Total O	0	0	
	11	99	35 35	0	U	
3	L	4	Total O	0	0	
	L	-1	4 4	0	U	
3	M	25	Total O	0	0	
3	171	20	25 25	U	U	
3	N	6	Total O	0	0	
		0	6 6		0	
3	О	21	Total O	0	0	
) 	U	41	21 21	U	U	
3	Р	1	Total O	0	0	
3	Г	1	1 1	U	U	



3 Residue-property plots (i)

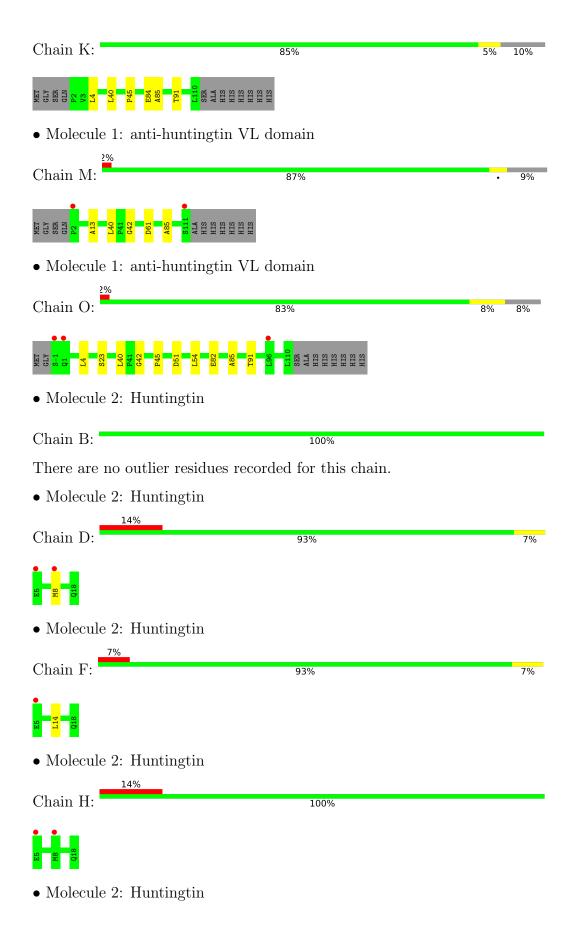
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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 (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: anti-huntingtin VL domain



• Molecule 1: anti-huntingtin VL domain







Chain J:	100%	
85 840 840 840 840 840 840 840 840 840 840		
• Molecule 2: Huntingtin		
Chain L:	100%	
2 2 3 3 3 3 3 3 3 3 3 3 3 3 3 3 3 3 3 3		
• Molecule 2: Huntingtin		
Chain N:	93%	7%
8 - 2 - 8		
• Molecule 2: Huntingtin		
Chain P:	100%	

There are no outlier residues recorded for this chain.



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants	53.09Å 89.66Å 95.00Å	Donositor
a, b, c, α , β , γ	90.00° 99.45° 90.00°	Depositor
Resolution (Å)	19.86 - 2.60	Depositor
rtesolution (A)	19.86 - 2.60	EDS
% Data completeness	88.9 (19.86-2.60)	Depositor
(in resolution range)	88.9 (19.86-2.60)	EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	3.88 (at 2.59Å)	Xtriage
Refinement program	REFMAC 5.5.0072	Depositor
D D.	0.216 , 0.278	Depositor
R, R_{free}	0.214 , 0.277	DCC
R_{free} test set	1221 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	15.0	Xtriage
Anisotropy	0.072	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.35, 24.4	EDS
L-test for twinning ²	$ < L > = 0.46, < L^2> = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	7652	wwPDB-VP
Average B, all atoms (Å ²)	23.0	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 21.00 % 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 7.6988e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.

²Theoretical values of <|L|>, $<L^2>$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.



¹Intensities estimated from amplitudes.

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).

Mal	Mol Chain		lengths	Bond	Bond angles		
IVIOI	Chain	RMSZ	# Z > 5	RMSZ	# Z > 5		
1	A	0.31	0/823	0.47	0/1124		
1	С	0.31	0/823	0.47	0/1124		
1	Е	0.32	0/844	0.46	0/1153		
1	G	0.31	0/823	0.46	0/1124		
1	I	0.31	0/829	0.47	0/1132		
1	K	0.32	0/823	0.45	0/1124		
1	M	0.33	0/829	0.47	0/1132		
1	О	0.31	0/838	0.44	0/1145		
2	В	0.40	0/119	0.37	0/153		
2	D	0.36	0/119	0.39	0/153		
2	F	0.40	0/119	0.37	0/153		
2	Н	0.36	0/119	0.38	0/153		
2	J	0.39	0/119	0.41	0/153		
2	L	0.37	0/119	0.37	0/153		
2	N	0.38	0/119	0.36	0/153		
2	Р	0.38	0/119	0.38	0/153		
All	All	0.32	0/7584	0.45	0/10282		

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	807	0	773	2	0
1	С	807	0	773	4	0
1	Ε	828	0	793	4	0
1	G	807	0	773	4	0
1	Ι	813	0	778	6	0
1	K	807	0	773	3	0
1	M	813	0	778	4	0
1	О	822	0	788	6	0
2	В	118	0	122	0	0
2	D	118	0	122	0	0
2	F	118	0	122	1	0
2	Н	118	0	122	0	0
2	J	118	0	122	0	0
2	L	118	0	122	0	0
2	N	118	0	122	0	0
2	Р	118	0	122	0	0
3	A	21	0	0	0	0
3	В	4	0	0	0	0
3	С	16	0	0	0	0
3	D	2	0	0	0	0
3	Ε	20	0	0	0	0
3	F	1	0	0	0	0
3	G	24	0	0	0	0
3	Н	1	0	0	0	0
3	I	22	0	0	0	0
3	J	1	0	0	0	0
3	K	35	0	0	0	0
3	L	4	0	0	0	0
3	M	25	0	0	0	0
3	N	6	0	0	0	0
3	О	21	0	0	0	0
3	Р	1	0	0	0	0
All	All	7652	0	7205	27	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 2.

All (27) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$\begin{array}{c} \text{Interatomic} \\ \text{distance (Å)} \end{array}$	$egin{aligned} \operatorname{Clash} \ \operatorname{overlap}\ (\mathring{\mathbf{A}}) \end{aligned}$
1:E:40:LEU:HD23	1:E:85:ALA:HB2	1.79	0.63
1:A:109:VAL:HG12	1:A:110:LEU:H	1.66	0.61



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A		Interatomic	Clash
Atom-1	Atom-2	${\rm distance}(\mathring{\rm A})$	overlap (Å)
1:C:42:GLY:HA2	1:E:45:PRO:HG3	1.85	0.59
1:I:45:PRO:HG3	1:O:42:GLY:HA2	1.84	0.58
1:M:40:LEU:HD23	1:M:85:ALA:HB2	1.85	0.58
1:G:36:TRP:HB2	1:G:49:MET:HB3	1.87	0.55
1:G:38:GLN:HB2	1:G:48:LEU:HD22	1.90	0.54
1:E:26:ASN:HA	1:E:30:GLY:HA3	1.89	0.53
1:I:40:LEU:HD23	1:I:85:ALA:HB2	1.91	0.53
1:O:40:LEU:HD23	1:O:85:ALA:HB2	1.90	0.52
1:I:84:GLU:HG2	1:I:108:THR:HA	1.91	0.51
1:C:57:PRO:HG3	1:M:61:ASP:HB2	1.94	0.50
1:A:109:VAL:HG12	1:A:110:LEU:N	2.26	0.49
1:C:40:LEU:HD23	1:C:85:ALA:HB2	1.94	0.49
1:I:42:GLY:HA2	1:O:45:PRO:HG3	1.94	0.48
1:I:12:ALA:HB3	1:I:18:VAL:HG12	1.96	0.48
1:O:4:LEU:HD11	1:O:91:THR:HG22	1.96	0.48
1:C:26:ASN:HA	1:C:30:GLY:HA3	1.97	0.46
1:K:40:LEU:HD23	1:K:85:ALA:HB2	1.98	0.46
1:O:51:ASP:HB3	1:O:54:LEU:HD12	1.99	0.44
1:M:13:ALA:HB2	1:O:23:SER:HB3	2.00	0.43
1:G:40:LEU:HD23	1:G:85:ALA:HB2	2.01	0.42
1:G:4:LEU:HD11	1:G:91:THR:HG22	2.02	0.42
1:K:45:PRO:HG3	1:M:42:GLY:HA2	2.01	0.42
1:I:4:LEU:HD11	1:I:91:THR:HG22	2.02	0.41
1:E:37:ILE:HG12	2:F:14:LEU:HD12	2.03	0.40
1:K:4:LEU:HD11	1:K:91:THR:HG22	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	107/121 (88%)	99 (92%)	7 (6%)	1 (1%)	17 35



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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Perce	\mathbf{ntiles}
1	\mathbf{C}	$107/121\ (88\%)$	103 (96%)	4 (4%)	0	100	100
1	${ m E}$	$110/121\ (91\%)$	106 (96%)	4 (4%)	0	100	100
1	G	$107/121\ (88\%)$	102 (95%)	5 (5%)	0	100	100
1	I	$108/121\ (89\%)$	102 (94%)	6 (6%)	0	100	100
1	K	107/121 (88%)	104 (97%)	3 (3%)	0	100	100
1	M	108/121 (89%)	105 (97%)	3 (3%)	0	100	100
1	О	$109/121\ (90\%)$	104 (95%)	5 (5%)	0	100	100
2	В	12/14~(86%)	12 (100%)	0	0	100	100
2	D	12/14~(86%)	12 (100%)	0	0	100	100
2	F	12/14~(86%)	12 (100%)	0	0	100	100
2	Н	12/14~(86%)	12 (100%)	0	0	100	100
2	J	12/14~(86%)	12 (100%)	0	0	100	100
2	L	12/14 (86%)	12 (100%)	0	0	100	100
2	N	12/14 (86%)	12 (100%)	0	0	100	100
2	Р	12/14 (86%)	12 (100%)	0	0	100	100
All	All	959/1080 (89%)	921 (96%)	37 (4%)	1 (0%)	51	75

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	84	GLU

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	90/100 (90%)	89 (99%)	1 (1%)	73 88
1	C	90/100 (90%)	89 (99%)	1 (1%)	73 88
1	E	93/100 (93%)	93 (100%)	0	100 100



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Mol	Chain	Analysed	Rotameric	Outliers	Perce	ntiles
1	G	90/100 (90%)	90 (100%)	0	100	100
1	I	91/100 (91%)	90 (99%)	1 (1%)	73	88
1	K	90/100 (90%)	89 (99%)	1 (1%)	73	88
1	M	91/100 (91%)	91 (100%)	0	100	100
1	О	92/100 (92%)	91 (99%)	1 (1%)	73	88
2	В	13/13 (100%)	13 (100%)	0	100	100
2	D	13/13 (100%)	12 (92%)	1 (8%)	13	25
2	F	13/13 (100%)	13 (100%)	0	100	100
2	Н	13/13 (100%)	13 (100%)	0	100	100
2	J	13/13 (100%)	13 (100%)	0	100	100
2	L	13/13 (100%)	13 (100%)	0	100	100
2	N	13/13 (100%)	12 (92%)	1 (8%)	13	25
2	Р	13/13 (100%)	13 (100%)	0	100	100
All	All	831/904 (92%)	824 (99%)	7 (1%)	81	92

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	86	ASP
1	С	18	VAL
2	D	8	MET
1	I	18	VAL
1	K	84	GLU
2	N	14	LEU
1	О	82	GLU

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

Mol	Chain	Res	Type
1	M	80	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 (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, 95^{th} 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>	# RSRZ > 2	$\mathbf{OWAB}(\mathbf{\mathring{A}}^2)$	Q<0.9
1	A	109/121 (90%)	-0.29	1 (0%) 84 82	16, 20, 28, 31	0
1	С	109/121 (90%)	-0.12	0 100 100	17, 24, 34, 37	0
1	Е	112/121 (92%)	-0.17	2 (1%) 68 64	16, 23, 35, 41	0
1	G	109/121 (90%)	-0.22	1 (0%) 84 82	16, 21, 34, 39	0
1	I	110/121 (90%)	-0.36	0 100 100	15, 18, 24, 27	0
1	K	109/121 (90%)	-0.34	0 100 100	15, 19, 27, 32	0
1	M	110/121 (90%)	-0.22	2 (1%) 68 64	15, 19, 27, 33	0
1	О	111/121 (91%)	-0.09	3 (2%) 54 48	15, 25, 40, 48	0
2	В	14/14 (100%)	-0.30	0 100 100	19, 24, 31, 40	0
2	D	14/14 (100%)	0.54	2 (14%) 2 1	30, 36, 52, 66	0
2	F	14/14 (100%)	0.14	1 (7%) 16 11	24, 31, 55, 69	0
2	Н	14/14 (100%)	0.29	2 (14%) 2 1	31, 38, 47, 63	0
2	J	14/14 (100%)	-0.05	1 (7%) 16 11	22, 26, 40, 45	0
2	L	14/14 (100%)	0.03	1 (7%) 16 11	21, 27, 43, 60	0
2	N	14/14 (100%)	0.14	1 (7%) 16 11	20, 25, 43, 62	0
2	Р	14/14 (100%)	0.34	0 100 100	40, 50, 60, 68	0
All	All	991/1080 (91%)	-0.19	17 (1%) 70 66	15, 22, 39, 69	0

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	N	5	GLU	5.4
1	M	2	PRO	4.1
1	O	-1	SER	3.2
2	D	8	MET	3.1
2	F	5	GLU	3.0



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Mol	Chain	Res	Type	RSRZ
1	M	111	SER	3.0
2	L	5	GLU	3.0
2	Н	5	GLU	2.8
2	Н	8	MET	2.7
1	Е	-1	SER	2.7
2	D	5	GLU	2.6
1	О	1	GLN	2.5
1	A	110	LEU	2.5
1	G	2	PRO	2.3
1	Е	26	ASN	2.2
1	О	96	LEU	2.0
2	J	5	GLU	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.

