



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 19, 2016 – 10:29 PM GMT

PDB ID : 5D68
Title : Crystal structure of KRIT1 ARD-FERM
Authors : Zhang, R.; Li, X.; Boggon, T.J.
Deposited on : 2015-08-11
Resolution : 2.91 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
<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 : unknown
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026982
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) : rb-20026982

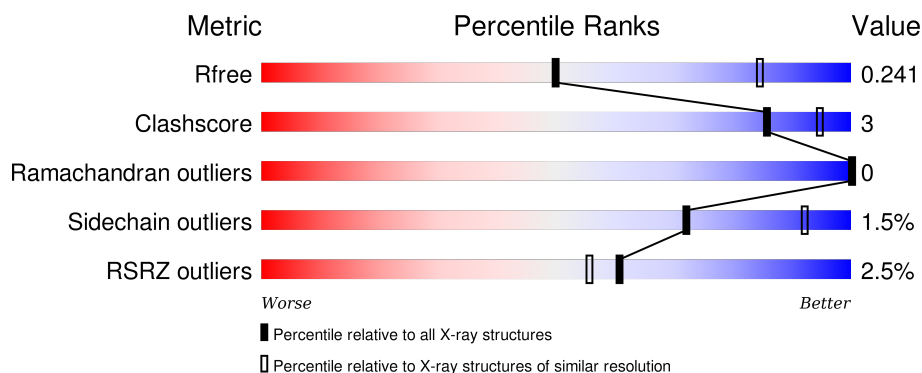
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

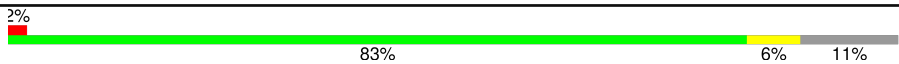
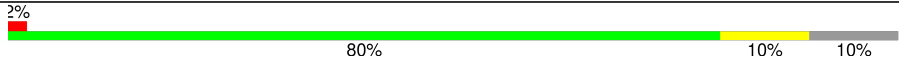

The reported resolution of this entry is 2.91 Å.

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	1451 (2.90-2.90)
Clashscore	102246	1668 (2.90-2.90)
Ramachandran outliers	100387	1630 (2.90-2.90)
Sidechain outliers	100360	1632 (2.90-2.90)
RSRZ outliers	91569	1456 (2.90-2.90)

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	486	
1	B	486	
1	C	486	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 10652 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 Krev interaction trapped protein 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	433	Total	C	N	O	S	0	0	0
			3524	2254	613	638	19			
1	B	437	Total	C	N	O	S	0	0	0
			3555	2269	619	648	19			
1	C	431	Total	C	N	O	S	0	0	0
			3507	2243	610	636	18			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	251	GLY	-	expression tag	UNP O00522
A	252	SER	-	expression tag	UNP O00522
A	253	ASP	-	expression tag	UNP O00522
A	254	ILE	-	expression tag	UNP O00522
A	255	LEU	-	expression tag	UNP O00522
A	256	GLN	-	expression tag	UNP O00522
A	257	GLY	-	expression tag	UNP O00522
A	258	THR	-	expression tag	UNP O00522
B	251	GLY	-	expression tag	UNP O00522
B	252	SER	-	expression tag	UNP O00522
B	253	ASP	-	expression tag	UNP O00522
B	254	ILE	-	expression tag	UNP O00522
B	255	LEU	-	expression tag	UNP O00522
B	256	GLN	-	expression tag	UNP O00522
B	257	GLY	-	expression tag	UNP O00522
B	258	THR	-	expression tag	UNP O00522
C	251	GLY	-	expression tag	UNP O00522
C	252	SER	-	expression tag	UNP O00522
C	253	ASP	-	expression tag	UNP O00522
C	254	ILE	-	expression tag	UNP O00522
C	255	LEU	-	expression tag	UNP O00522
C	256	GLN	-	expression tag	UNP O00522
C	257	GLY	-	expression tag	UNP O00522

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Chain	Residue	Modelled	Actual	Comment	Reference
C	258	THR	-	expression tag	UNP O00522

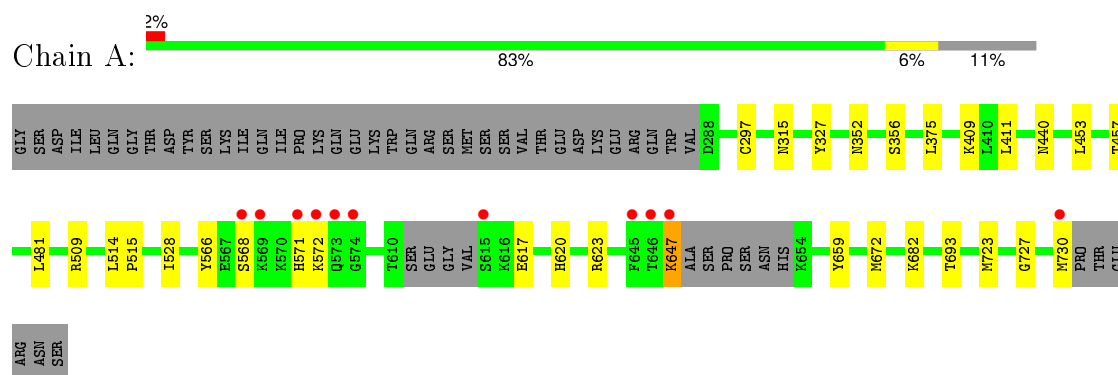
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	19	Total	O	0	0
			19	19		
2	B	26	Total	O	0	0
			26	26		
2	C	21	Total	O	0	0
			21	21		

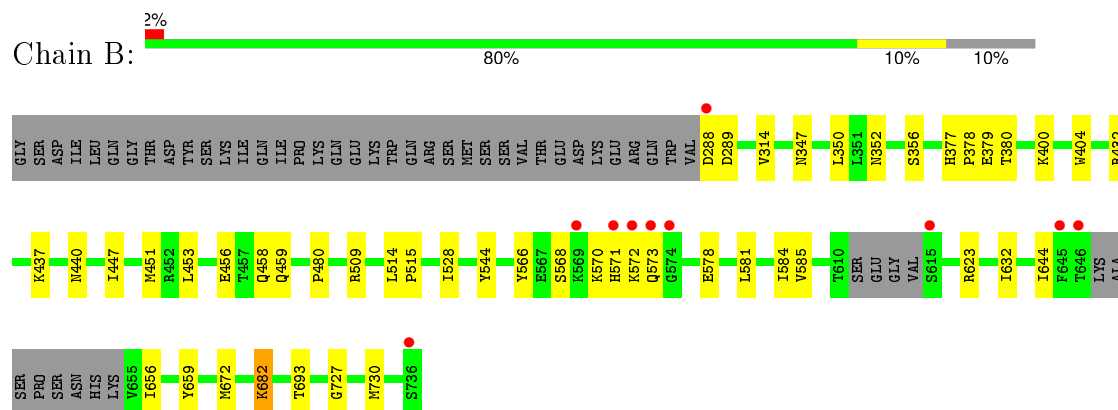
3 Residue-property plots [i](#)

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.

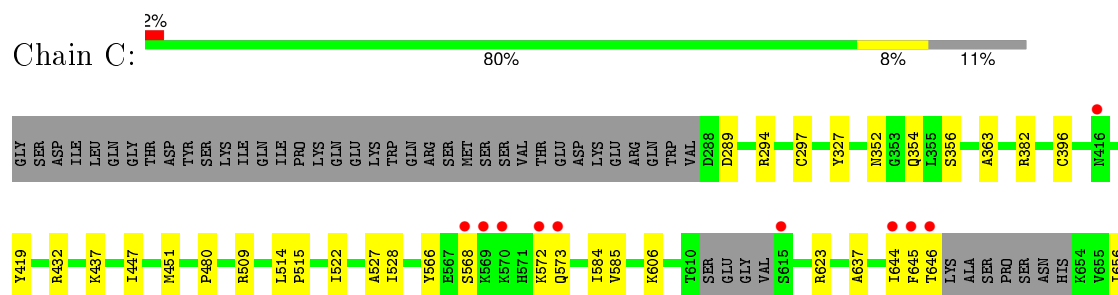
• Molecule 1: Krev interaction trapped protein 1



• Molecule 1: Krev interaction trapped protein 1



• Molecule 1: Krev interaction trapped protein 1





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	125.73Å 80.24Å 213.15Å 90.00° 91.69° 90.00°	Depositor
Resolution (Å)	49.37 – 2.91 49.37 – 2.91	Depositor EDS
% Data completeness (in resolution range)	99.6 (49.37-2.91) 90.3 (49.37-2.91)	Depositor EDS
R_{merge}	0.19	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.77 (at 2.91Å)	Xtriage
Refinement program	PHENIX (phenix.refine: dev_1760)	Depositor
R, R_{free}	0.207 , 0.246 0.205 , 0.241	Depositor DCC
R_{free} test set	2156 reflections (5.08%)	DCC
Wilson B-factor (Å ²)	37.6	Xtriage
Anisotropy	0.629	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 30.1	EDS
Estimated twinning fraction	0.018 for -h,-k,l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 46891 reflections	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	10652	wwPDB-VP
Average B, all atoms (Å ²)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.31% 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	0.27	0/3604	0.42	0/4872
1	B	0.26	0/3636	0.42	0/4917
1	C	0.26	0/3587	0.43	0/4851
All	All	0.26	0/10827	0.42	0/14640

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	3524	0	3540	13	0
1	B	3555	0	3558	24	0
1	C	3507	0	3518	23	0
2	A	19	0	0	0	0
2	B	26	0	0	0	0
2	C	21	0	0	1	0
All	All	10652	0	10616	59	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 3.

All (59) 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:B:451:MET:HB3	1:B:453:LEU:HD13	1.77	0.67
1:C:645:PHE:N	1:C:646:THR:HA	2.17	0.59
1:C:297:CYS:HB3	1:C:327:TYR:HB3	1.85	0.58
1:B:727:GLY:HA2	1:B:730:MET:HG3	1.86	0.58
1:A:297:CYS:HB3	1:A:327:TYR:HB3	1.90	0.53
1:C:352:ASN:ND2	1:C:354:GLN:OE1	2.43	0.51
1:C:352:ASN:HB3	1:C:356:SER:H	1.76	0.51
1:B:566:TYR:CE2	1:B:571:HIS:HB2	2.45	0.51
1:C:572:LYS:HG2	1:C:573:GLN:HG3	1.93	0.51
1:A:352:ASN:HB3	1:A:356:SER:H	1.75	0.50
1:A:659:TYR:CE2	1:A:672:MET:HG3	2.46	0.50
1:C:528:ILE:HD12	1:C:623:ARG:HG3	1.94	0.50
1:B:447:ILE:HD12	1:B:480:PRO:HG2	1.93	0.49
1:C:568:SER:O	1:C:572:LYS:HB2	2.14	0.48
1:B:572:LYS:HG2	1:B:573:GLN:HG3	1.96	0.47
1:B:437:LYS:HE3	1:C:702:MET:O	2.14	0.47
1:C:419:TYR:CD1	1:C:437:LYS:HA	2.50	0.47
1:C:584:ILE:HG13	1:C:585:VAL:HG13	1.97	0.46
1:A:568:SER:O	1:A:572:LYS:HB2	2.16	0.46
1:C:447:ILE:HD12	1:C:480:PRO:HG2	1.97	0.46
1:A:566:TYR:CE2	1:A:571:HIS:HB2	2.50	0.46
1:B:288:ASP:OD1	1:B:289:ASP:N	2.49	0.46
1:A:727:GLY:O	1:A:730:MET:HG3	2.16	0.46
1:B:352:ASN:HB3	1:B:356:SER:H	1.81	0.45
1:C:644:ILE:HG12	1:C:656:ILE:HB	1.98	0.45
1:C:682:LYS:HB2	1:C:682:LYS:HE3	1.81	0.45
1:A:617:GLU:HB2	1:A:620:HIS:ND1	2.32	0.45
1:B:544:TYR:HB3	1:B:632:ILE:HD13	1.99	0.45
1:C:382:ARG:HG2	2:C:809:HOH:O	2.17	0.44
1:B:568:SER:O	1:B:572:LYS:HB2	2.16	0.44
1:A:528:ILE:HD12	1:A:623:ARG:HG3	2.00	0.44
1:B:528:ILE:HD12	1:B:623:ARG:HG3	1.98	0.44
1:C:659:TYR:HB2	1:C:670:LEU:HB2	1.99	0.43
1:B:347:ASN:HB3	1:B:350:LEU:HB2	1.99	0.43
1:B:377:HIS:HA	1:B:378:PRO:HD3	1.87	0.43
1:A:440:ASN:HA	1:A:481:LEU:HD12	2.00	0.43
1:C:566:TYR:CD1	1:C:606:LYS:HG2	2.53	0.43
1:C:363:ALA:HB1	1:C:396:CYS:HB2	2.00	0.43
1:C:637:ALA:HA	1:C:663:ASN:HB3	2.00	0.43
1:A:375:LEU:HD11	1:A:411:LEU:HD23	2.00	0.42
1:C:514:LEU:HA	1:C:515:PRO:HD3	1.90	0.42
1:B:400:LYS:HA	1:B:404:TRP:CG	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:514:LEU:HD12	1:A:515:PRO:HD2	2.02	0.42
1:C:289:ASP:OD1	1:C:294:ARG:NH1	2.52	0.42
1:B:514:LEU:HA	1:B:515:PRO:HD3	1.94	0.42
1:C:514:LEU:HD12	1:C:515:PRO:HD2	2.02	0.42
1:B:456:GLU:O	1:B:459:GLN:HG2	2.20	0.42
1:B:659:TYR:CE2	1:B:672:MET:HG3	2.54	0.42
1:B:379:GLU:HB3	1:B:440:ASN:ND2	2.35	0.42
1:C:432:ARG:HD2	1:C:451:MET:HE3	2.02	0.41
1:B:682:LYS:HE3	1:B:682:LYS:HB2	1.78	0.41
1:B:644:ILE:HG12	1:B:656:ILE:HB	2.02	0.41
1:A:453:LEU:HB3	1:A:457:THR:HG23	2.02	0.41
1:B:377:HIS:HB3	1:B:380:THR:OG1	2.20	0.41
1:B:581:LEU:O	1:B:585:VAL:HG22	2.20	0.41
1:B:570:LYS:HE3	1:B:571:HIS:CE1	2.55	0.41
1:B:584:ILE:HG13	1:B:585:VAL:HG13	2.02	0.41
1:A:647:LYS:NZ	1:A:647:LYS:HB2	2.35	0.40
1:C:522:ILE:HD11	1:C:527:ALA:CB	2.51	0.40

There are no symmetry-related clashes.

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	427/486 (88%)	416 (97%)	11 (3%)	0	100	100
1	B	431/486 (89%)	418 (97%)	13 (3%)	0	100	100
1	C	425/486 (87%)	412 (97%)	13 (3%)	0	100	100
All	All	1283/1458 (88%)	1246 (97%)	37 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	390/439 (89%)	383 (98%)	7 (2%)	66	90
1	B	394/439 (90%)	387 (98%)	7 (2%)	66	90
1	C	388/439 (88%)	384 (99%)	4 (1%)	82	95
All	All	1172/1317 (89%)	1154 (98%)	18 (2%)	72	92

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	315	ASN
1	A	409	LYS
1	A	509	ARG
1	A	647	LYS
1	A	682	LYS
1	A	693	THR
1	A	723	MET
1	B	314	VAL
1	B	432	ARG
1	B	458	GLN
1	B	509	ARG
1	B	578	GLU
1	B	682	LYS
1	B	693	THR
1	C	509	ARG
1	C	682	LYS
1	C	693	THR
1	C	729	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	689	GLN
1	C	726	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	433/486 (89%)	-0.06	11 (2%) 61 55	21, 40, 83, 128	0
1	B	437/486 (89%)	-0.09	10 (2%) 64 59	23, 42, 87, 125	0
1	C	431/486 (88%)	-0.05	12 (2%) 56 50	23, 41, 83, 134	0
All	All	1301/1458 (89%)	-0.07	33 (2%) 61 55	21, 41, 84, 134	0

All (33) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	646	THR	7.1
1	A	572	LYS	5.6
1	A	573	GLN	5.0
1	B	573	GLN	4.1
1	A	571	HIS	4.1
1	A	574	GLY	3.9
1	A	569	LYS	3.3
1	C	644	ILE	3.3
1	B	736	SER	3.3
1	C	572	LYS	3.2
1	A	730	MET	3.2
1	B	574	GLY	3.1
1	B	646	THR	2.9
1	C	729	LEU	2.9
1	C	416	ASN	2.9
1	B	288	ASP	2.8
1	A	615	SER	2.8
1	B	615	SER	2.7
1	C	645	PHE	2.7
1	B	571	HIS	2.7
1	A	645	PHE	2.6
1	A	568	SER	2.6
1	A	646	THR	2.6

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Mol	Chain	Res	Type	RSRZ
1	C	569	LYS	2.5
1	C	615	SER	2.5
1	C	570	LYS	2.5
1	B	569	LYS	2.5
1	B	572	LYS	2.5
1	C	568	SER	2.5
1	C	573	GLN	2.4
1	C	727	GLY	2.2
1	B	645	PHE	2.2
1	A	647	LYS	2.1

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.