



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 31, 2021 – 12:04 PM EDT

PDB ID : 7K28
Title : Kelch domain of human KEAP1 bound to Nrf2 peptide, ADEETGEFL
Authors : Muellers, S.N.; Allen, K.N.
Deposited on : 2020-09-08
Resolution : 2.15 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Xtriage (Phenix) : 1.13
EDS : 2.18
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)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.18

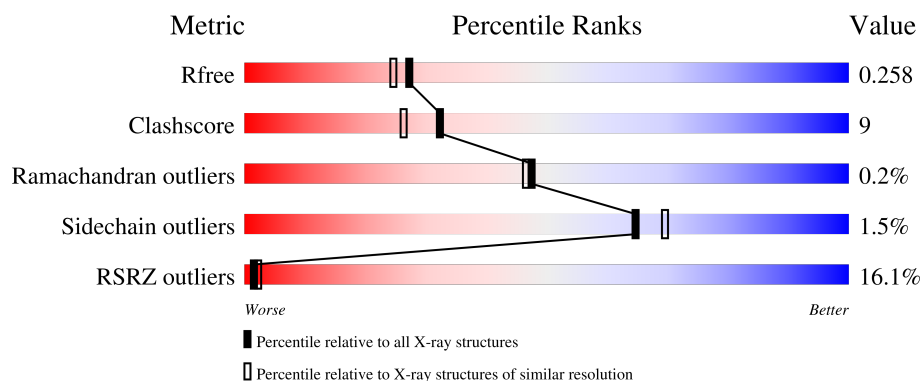
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.15 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	1479 (2.16-2.16)
Clashscore	141614	1585 (2.16-2.16)
Ramachandran outliers	138981	1560 (2.16-2.16)
Sidechain outliers	138945	1559 (2.16-2.16)
RSRZ outliers	127900	1456 (2.16-2.16)

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	290	<div> <div>9%</div> <div>83%</div> <div>14%</div> <div>..</div> </div>
1	B	290	<div> <div>23%</div> <div>84%</div> <div>13%</div> <div>.</div> </div>
2	P	11	<div> <div>73%</div> <div>9%</div> <div>18%</div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4424 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 Kelch-like ECH-associated protein 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	284	Total	C	N	O	S	0	1	0
			2175	1355	393	411	16			
1	B	284	Total	C	N	O	S	0	1	0
			2179	1356	394	413	16			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	540	ALA	GLU	conflict	UNP Q14145
A	542	ALA	GLU	conflict	UNP Q14145
A	613	SER	CYS	conflict	UNP Q14145
B	540	ALA	GLU	conflict	UNP Q14145
B	542	ALA	GLU	conflict	UNP Q14145
B	613	SER	CYS	conflict	UNP Q14145

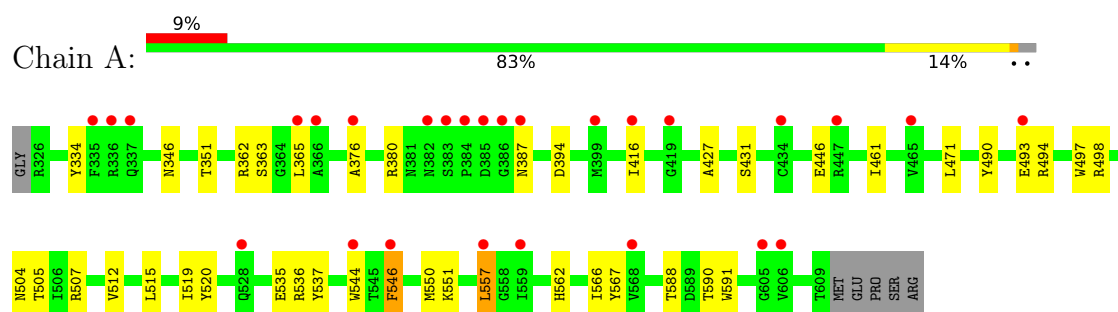
- Molecule 2 is a protein called Nrf2 peptide,ADEETGEFL.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	P	9	Total	C	N	O	0	0	0
			70	43	9	18			

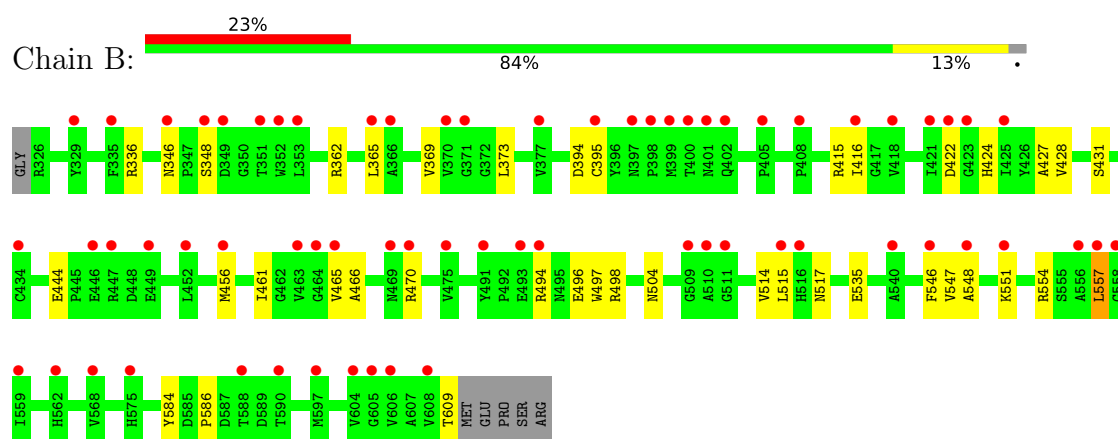
3 Residue-property plots

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.

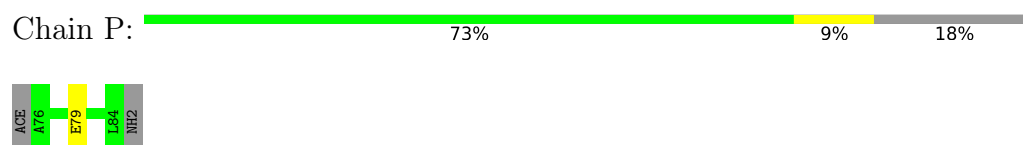
• Molecule 1: Kelch-like ECH-associated protein 1



• Molecule 1: Kelch-like ECH-associated protein 1



• Molecule 2: Nrf2 peptide, ADEETGEFL



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	162.01Å 68.61Å 77.35Å 90.00° 117.90° 90.00°	Depositor
Resolution (Å)	29.46 – 2.15 29.46 – 2.15	Depositor EDS
% Data completeness (in resolution range)	99.4 (29.46-2.15) 99.4 (29.46-2.15)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.71 (at 2.16Å)	Xtriage
Refinement program	PHENIX 1.17.1_3660	Depositor
R, R_{free}	0.238 , 0.258 0.238 , 0.258	Depositor DCC
R_{free} test set	2007 reflections (4.91%)	wwPDB-VP
Wilson B-factor (Å ²)	36.9	Xtriage
Anisotropy	0.315	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 35.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.010 for -h-2*k,l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	4424	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 5.78% 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	0.29	0/2231	0.56	0/3040
1	B	0.28	0/2234	0.53	0/3043
2	P	0.28	0/70	0.53	0/93
All	All	0.28	0/4535	0.55	0/6176

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	2175	0	2071	44	0
1	B	2179	0	2079	33	0
2	P	70	0	56	1	0
All	All	4424	0	4206	76	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 9.

All (76) 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:A:535:GLU:HB3	1:A:546:PHE:CD1	1.51	1.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:535:GLU:HB3	1:A:546:PHE:CE1	1.85	1.11
1:B:494:ARG:HG3	1:B:496:GLU:HG2	1.46	0.98
1:A:535:GLU:HB3	1:A:546:PHE:HD1	1.27	0.96
1:A:535:GLU:CB	1:A:546:PHE:CD1	2.47	0.96
1:B:504:ASN:HB2	1:B:546:PHE:CZ	2.04	0.92
1:A:537:TYR:CD1	1:A:544:TRP:CD1	2.61	0.89
1:A:537:TYR:HD1	1:A:544:TRP:CD1	1.90	0.88
1:B:466:ALA:HB1	1:B:514:VAL:HG23	1.58	0.84
1:A:537:TYR:CD1	1:A:544:TRP:NE1	2.46	0.83
1:A:535:GLU:CB	1:A:546:PHE:CE1	2.62	0.82
1:B:428:VAL:HG23	1:B:456:MET:SD	2.28	0.74
1:B:428:VAL:CG2	1:B:456:MET:SD	2.78	0.72
1:B:504:ASN:HB2	1:B:546:PHE:CE2	2.26	0.71
1:A:537:TYR:HB2	1:A:544:TRP:CE2	2.28	0.69
1:A:505:THR:HG1	1:A:546:PHE:HZ	1.42	0.68
1:A:535:GLU:CG	1:A:546:PHE:CE1	2.77	0.68
1:B:504:ASN:CB	1:B:546:PHE:CE2	2.77	0.68
1:A:535:GLU:CD	1:A:546:PHE:CE1	2.70	0.64
1:B:415:ARG:NH2	2:P:79:GLU:OE1	2.25	0.62
1:B:416:ILE:HD11	1:B:427:ALA:HB1	1.81	0.62
1:A:537:TYR:HB2	1:A:544:TRP:CD2	2.35	0.62
1:A:535:GLU:CG	1:A:546:PHE:HE1	2.14	0.61
1:B:504:ASN:CB	1:B:546:PHE:CZ	2.82	0.61
1:A:515:LEU:HD22	1:A:566:ILE:HG13	1.82	0.60
1:B:456:MET:HG2	1:B:497:TRP:CE2	2.36	0.60
1:B:428:VAL:HG11	1:B:465:VAL:HG21	1.86	0.58
1:A:416:ILE:HD11	1:A:427:ALA:HB1	1.85	0.57
1:A:550:MET:HG2	1:A:591:TRP:CE2	2.40	0.57
1:B:362:ARG:NH1	1:B:394:ASP:OD2	2.39	0.55
1:A:493:GLU:HG3	1:A:494:ARG:HG3	1.89	0.55
1:B:428:VAL:CG1	1:B:465:VAL:HG21	2.37	0.55
1:B:365:LEU:H	1:B:365:LEU:HD23	1.72	0.54
1:B:456:MET:HG2	1:B:497:TRP:CZ2	2.43	0.53
1:A:380:ARG:NH1	1:A:387:ASN:HB3	2.24	0.53
1:B:373:LEU:HB3	1:B:395:CYS:SG	2.49	0.53
1:B:362:ARG:HG3	1:B:365:LEU:HD13	1.89	0.53
1:A:519:ILE:O	1:A:536:ARG:HA	2.09	0.52
1:A:535:GLU:CB	1:A:546:PHE:HD1	2.06	0.51
1:B:346:ASN:HD21	1:B:348:SER:HB2	1.75	0.51
1:B:428:VAL:HG21	1:B:456:MET:SD	2.52	0.49
1:A:380:ARG:HH12	1:A:387:ASN:HB3	1.78	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:504:ASN:ND2	1:A:544:TRP:O	2.42	0.47
1:A:334:TYR:C	1:A:334:TYR:CD1	2.88	0.47
1:A:346:ASN:HB3	1:A:351:THR:HG22	1.96	0.47
1:B:415:ARG:HD3	1:B:461:ILE:HD12	1.96	0.47
1:A:588:THR:O	1:A:590:THR:HG23	2.16	0.46
1:B:369:VAL:HG21	1:B:609:THR:HB	1.98	0.46
1:B:346:ASN:ND2	1:B:348:SER:HB2	2.30	0.45
1:B:548:ALA:HB3	1:B:584:TYR:HE1	1.81	0.45
1:A:512:VAL:HA	1:A:520:TYR:O	2.17	0.45
1:A:446:GLU:H	1:A:446:GLU:CD	2.18	0.45
1:B:424:HIS:CE1	1:B:444:GLU:HG2	2.51	0.45
1:A:471:LEU:HD11	1:B:517:ASN:HB2	1.99	0.45
1:B:431:SER:HB3	1:B:461:ILE:HG21	1.99	0.44
1:A:334:TYR:CD1	1:A:334:TYR:O	2.70	0.44
1:A:334:TYR:HB2	1:A:363:SER:HB3	1.98	0.44
1:A:334:TYR:HB2	1:A:363:SER:CB	2.47	0.44
1:A:362:ARG:NH1	1:A:394:ASP:OD1	2.50	0.44
1:A:365:LEU:HD23	1:A:365:LEU:H	1.81	0.44
1:A:535:GLU:HG2	1:A:546:PHE:HE1	1.83	0.43
1:A:537:TYR:CD1	1:A:544:TRP:CE2	3.06	0.43
1:A:507:ARG:HH21	1:A:546:PHE:HE1	1.66	0.43
1:A:557:LEU:HD23	1:A:557:LEU:H	1.84	0.42
1:A:365:LEU:HD12	1:A:376:ALA:HB1	2.01	0.42
1:A:431:SER:HB3	1:A:461:ILE:HG21	2.01	0.42
1:B:535:GLU:HB3	1:B:546:PHE:CD1	2.54	0.42
1:B:554:ARG:HG3	1:B:557:LEU:HD22	2.02	0.41
1:A:546:PHE:HD1	1:A:546:PHE:HA	1.77	0.41
1:A:334:TYR:O	1:A:334:TYR:HD1	2.02	0.41
1:B:498:ARG:HD2	1:B:498:ARG:HA	2.01	0.41
1:A:490:TYR:HB2	1:A:497:TRP:CE2	2.56	0.41
1:B:515:LEU:HD21	1:B:586:PRO:HG3	2.02	0.41
1:B:422:ASP:OD1	1:B:470:ARG:CZ	2.69	0.41
1:B:557:LEU:H	1:B:557:LEU:HD23	1.86	0.40
1:A:562:HIS:HB3	1:A:567:TYR:CE2	2.57	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	283/290 (98%)	272 (96%)	11 (4%)	0	100	100
1	B	283/290 (98%)	272 (96%)	10 (4%)	1 (0%)	34	29
2	P	7/11 (64%)	7 (100%)	0	0	100	100
All	All	573/591 (97%)	551 (96%)	21 (4%)	1 (0%)	47	46

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	547	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	228/234 (97%)	224 (98%)	4 (2%)	59	63
1	B	229/234 (98%)	226 (99%)	3 (1%)	69	74
2	P	7/7 (100%)	7 (100%)	0	100	100
All	All	464/475 (98%)	457 (98%)	7 (2%)	65	69

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

Mol	Chain	Res	Type
1	A	498	ARG
1	A	546	PHE

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Mol	Chain	Res	Type
1	A	551	LYS
1	A	557	LEU
1	B	336	ARG
1	B	551	LYS
1	B	557	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	346	ASN
1	B	359	GLN
1	B	402	GLN
1	B	517	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 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	284/290 (97%)	0.54	27 (9%) 8 12	20, 29, 49, 107	0
1	B	284/290 (97%)	1.20	66 (23%) 0 0	26, 54, 92, 116	0
2	P	9/11 (81%)	0.11	0 100 100	31, 37, 46, 49	0
All	All	577/591 (97%)	0.86	93 (16%) 1 2	20, 38, 86, 116	0

All (93) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	399	MET	6.7
1	A	384	PRO	6.0
1	B	351	THR	6.0
1	B	371	GLY	4.6
1	B	335	PHE	4.6
1	B	446	GLU	4.3
1	B	402	GLN	4.3
1	A	546	PHE	4.2
1	A	399	MET	4.1
1	B	557	LEU	4.1
1	A	386	GLY	4.1
1	B	604	VAL	4.0
1	B	470	ARG	4.0
1	B	447	ARG	3.9
1	B	516	HIS	3.9
1	B	546	PHE	3.9
1	B	365	LEU	3.9
1	B	559	ILE	3.9
1	B	353	LEU	3.8
1	B	575	HIS	3.8
1	B	329	TYR	3.7
1	B	366	ALA	3.7
1	B	608	VAL	3.6

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Mol	Chain	Res	Type	RSRZ
1	B	493	GLU	3.6
1	B	494	ARG	3.6
1	B	416	ILE	3.5
1	A	385	ASP	3.5
1	A	366	ALA	3.5
1	B	422	ASP	3.4
1	B	400	THR	3.4
1	B	346	ASN	3.3
1	B	421	ILE	3.3
1	B	395	CYS	3.2
1	A	383	SER	3.1
1	B	423	GLY	3.0
1	B	606	VAL	3.0
1	B	418	VAL	3.0
1	B	465	VAL	3.0
1	A	416	ILE	2.9
1	A	434[A]	CYS	2.9
1	B	491	TYR	2.9
1	B	349	ASP	2.9
1	A	606	VAL	2.8
1	A	335	PHE	2.8
1	B	556	ALA	2.8
1	B	401	ASN	2.8
1	B	370	VAL	2.8
1	A	544	TRP	2.7
1	A	336	ARG	2.7
1	A	365	LEU	2.7
1	B	515	LEU	2.6
1	B	562	HIS	2.6
1	B	540	ALA	2.6
1	B	551	LYS	2.6
1	B	597	MET	2.5
1	B	475	VAL	2.5
1	B	588	THR	2.5
1	A	568	VAL	2.5
1	A	382	ASN	2.5
1	B	605	GLY	2.4
1	B	510	ALA	2.4
1	B	348	SER	2.4
1	B	568	VAL	2.4
1	B	397	ASN	2.4
1	B	408	PRO	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	376	ALA	2.4
1	B	464	GLY	2.4
1	A	605	GLY	2.4
1	A	557	LEU	2.3
1	B	352	TRP	2.3
1	A	387	ASN	2.3
1	B	398	PRO	2.3
1	B	377	VAL	2.3
1	A	559	ILE	2.3
1	A	528	GLN	2.2
1	B	452	LEU	2.2
1	B	469	ASN	2.2
1	A	465	VAL	2.2
1	B	463	VAL	2.2
1	B	449	GLU	2.2
1	B	405	PRO	2.1
1	B	511	GLY	2.1
1	B	558	GLY	2.1
1	B	509	GLY	2.1
1	A	493	GLU	2.1
1	B	456	MET	2.1
1	B	548	ALA	2.1
1	B	590	THR	2.1
1	A	419	GLY	2.0
1	B	425	ILE	2.0
1	B	434[A]	CYS	2.0
1	A	337	GLN	2.0
1	A	447	ARG	2.0

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 monosaccharides in this entry.

6.4 Ligands ⓘ

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.