



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 05:56 AM UTC

PDB ID : 8Y7F / pdb_00008y7f
Title : Crystal structure of CARF domain-truncated Csx1-Crn2 from *Marinitoga* sp.
Authors : Zhang, D.; Yuan, C.; Lin, Z.
Deposited on : 2024-02-04
Resolution : 2.95 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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

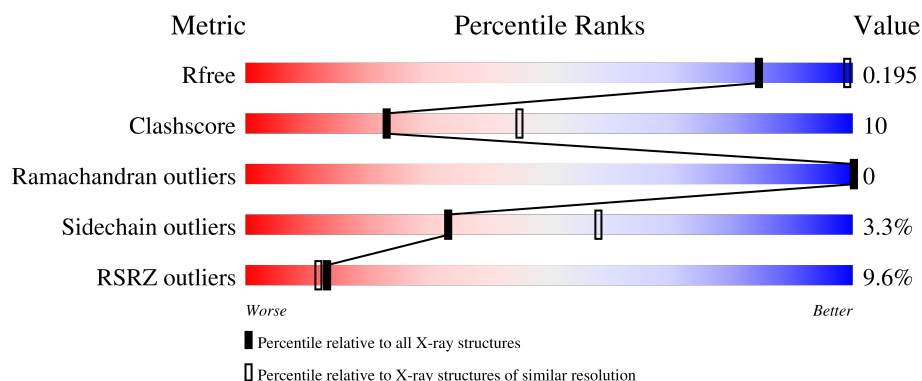
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.95 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}	180053	1130 (2.98-2.94)
Clashscore	190562	1157 (2.98-2.94)
Ramachandran outliers	187476	1101 (2.98-2.94)
Sidechain outliers	187428	1101 (2.98-2.94)
RSRZ outliers	180081	1130 (2.98-2.94)

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	382	7% 73% 23% .
1	B	382	11% 70% 23% 5%

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 6023 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 CRISPR-associated protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	368	Total	C	N	O	S	0	0	0
			3034	1980	488	555	11			
1	B	364	Total	C	N	O	S	0	0	0
			2989	1947	481	550	11			

4 Data and refinement statistics

Property	Value	Source
Space group	P 43 3 2	Depositor
Cell constants a, b, c, α , β , γ	210.58Å 210.58Å 210.58Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.53 – 2.95 40.53 – 2.95	Depositor EDS
% Data completeness (in resolution range)	100.0 (40.53-2.95) 99.9 (40.53-2.95)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.25 (at 2.95Å)	Xtriage
Refinement program	PHENIX (1.20.1_4487: ???)	Depositor
R, R_{free}	0.200 , 0.230 (Not available) , 0.195	Depositor DCC
R_{free} test set	1658 reflections (4.86%)	wwPDB-VP
Wilson B-factor (Å ²)	58.0	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 49.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	6023	wwPDB-VP
Average B, all atoms (Å ²)	60.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.77% 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

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	0.50	0/3083	0.69	5/4136 (0.1%)
1	B	0.49	0/3037	0.63	1/4076 (0.0%)
All	All	0.49	0/6120	0.66	6/8212 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	B	0	1
All	All	0	2

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	310	ASN	N-CA-CB	-7.01	98.97	110.40
1	A	481	GLU	CA-CB-CG	6.45	127.01	114.10
1	A	481	GLU	CA-C-N	-6.21	111.87	120.38
1	A	481	GLU	C-N-CA	-6.21	111.87	120.38
1	B	215	GLN	CB-CA-C	6.14	121.29	110.85
1	A	481	GLU	N-CA-CB	6.14	119.14	110.12

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	310	ASN	Sidechain
1	B	270	ARG	Sidechain

5.2 Too-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 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	3034	0	3146	66	1
1	B	2989	0	3088	66	1
All	All	6023	0	6234	126	1

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 10.

All (126) 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:256:ASN:HB2	1:A:310:ASN:HD21	1.31	0.95
1:A:530:LEU:HD11	1:A:545:MET:HE2	1.59	0.84
1:B:355:GLU:OE1	1:B:376:TYR:OH	1.99	0.80
1:B:376:TYR:HD2	1:B:391:GLU:HG2	1.54	0.72
1:A:256:ASN:CB	1:A:310:ASN:HD21	2.05	0.69
1:A:393:LEU:HA	1:A:429:MET:HE2	1.76	0.66
1:B:433:ILE:HG22	1:B:437:MET:HE2	1.77	0.66
1:B:376:TYR:CD2	1:B:391:GLU:HG2	2.31	0.65
1:A:256:ASN:HD22	1:A:310:ASN:ND2	1.95	0.65
1:A:195:LEU:HA	1:A:207:MET:CE	2.30	0.62
1:B:475:GLU:O	1:B:479:ILE:HG12	2.00	0.61
1:A:195:LEU:HD23	1:A:207:MET:HE1	1.82	0.60
1:B:193:ILE:HD11	1:B:206:ARG:NE	2.16	0.60
1:A:433:ILE:HG22	1:A:437:MET:HE2	1.84	0.60
1:B:384:LYS:NZ	1:B:385:GLU:H	1.99	0.60
1:A:445:LYS:HG2	1:A:446:LYS:H	1.67	0.58
1:B:482:ASN:OD1	1:B:485:LYS:HE2	2.04	0.58
1:A:201:TYR:HB2	1:A:203:ILE:HG12	1.86	0.58
1:A:332:ILE:HD11	1:A:423:LEU:HD13	1.86	0.57
1:B:433:ILE:CG2	1:B:437:MET:HE2	2.35	0.57
1:A:407:GLY:HA2	1:B:248:PRO:CB	2.34	0.57
1:B:351:ILE:HG23	1:B:437:MET:HE1	1.87	0.57
1:B:438:LYS:HG3	1:B:439:ASP:N	2.20	0.56
1:B:193:ILE:HD11	1:B:206:ARG:HE	1.71	0.56
1:A:250:PRO:HG2	1:A:253:LYS:HG3	1.88	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:441:LYS:HZ1	1:A:444:SER:N	2.05	0.55
1:B:449:LEU:HB2	1:B:489:LEU:HD11	1.87	0.55
1:A:513:LYS:HD3	1:A:515:LYS:HD3	1.89	0.55
1:A:256:ASN:HB2	1:A:310:ASN:ND2	2.13	0.54
1:B:352:LYS:HD3	1:B:376:TYR:CE2	2.42	0.54
1:A:322:GLU:HG2	1:A:434:LEU:HD13	1.90	0.54
1:A:527:VAL:HB	1:A:557:TYR:HB2	1.89	0.54
1:A:375:SER:HB3	1:A:391:GLU:OE2	2.08	0.54
1:B:345:ARG:HH22	1:B:369:GLU:CD	2.17	0.53
1:A:433:ILE:CG2	1:A:437:MET:HE2	2.39	0.53
1:A:481:GLU:O	1:A:482:ASN:C	2.47	0.53
1:B:500:GLU:HG2	1:B:539:ILE:CD1	2.40	0.52
1:A:251:PRO:HA	1:A:324:MET:HE3	1.91	0.52
1:B:258:ILE:HG21	1:B:304:PRO:O	2.09	0.52
1:B:348:LEU:HD11	1:B:397:VAL:HG21	1.91	0.52
1:B:355:GLU:HG2	1:B:355:GLU:O	2.07	0.51
1:A:194:ALA:C	1:A:207:MET:HE2	2.36	0.51
1:A:407:GLY:HA2	1:B:248:PRO:HB3	1.92	0.51
1:A:373:MET:HE2	1:A:374:TYR:CZ	2.45	0.51
1:B:384:LYS:O	1:B:388:LYS:HG2	2.11	0.51
1:B:384:LYS:HZ3	1:B:385:GLU:H	1.60	0.50
1:A:251:PRO:HB3	1:A:324:MET:HE3	1.94	0.50
1:B:500:GLU:HG2	1:B:539:ILE:HD12	1.93	0.49
1:A:547:PHE:O	1:A:550:GLU:CB	2.61	0.49
1:A:407:GLY:HA2	1:B:248:PRO:HB2	1.94	0.49
1:A:205:GLU:OE2	1:B:200:ARG:NH2	2.45	0.49
1:A:348:LEU:HD11	1:A:397:VAL:HG21	1.95	0.49
1:B:210:VAL:O	1:B:214:ILE:HG13	2.12	0.49
1:A:315:ILE:O	1:A:320:LYS:NZ	2.42	0.48
1:B:440:LEU:HD22	1:B:441:LYS:H	1.78	0.48
1:A:470:LYS:HE2	1:A:557:TYR:OH	2.13	0.48
1:B:260:LYS:O	1:B:264:ILE:HG13	2.13	0.48
1:A:248:PRO:HB2	1:B:407:GLY:HA2	1.94	0.48
1:A:393:LEU:CA	1:A:429:MET:HE2	2.43	0.48
1:B:349:ILE:HD11	1:B:372:TRP:HB2	1.95	0.47
1:B:297:LYS:HB2	1:B:297:LYS:HE2	1.41	0.47
1:A:265:LEU:O	1:A:269:ILE:HG12	2.14	0.47
1:A:281:ASN:O	1:A:285:ILE:HG13	2.13	0.47
1:A:498:VAL:HG22	1:A:545:MET:HE1	1.96	0.47
1:B:217:ILE:O	1:B:220:LYS:HG2	2.15	0.47
1:A:345:ARG:NH2	1:A:369:GLU:OE2	2.47	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:446:LYS:HD3	1:B:448:TYR:CE1	2.50	0.47
1:A:256:ASN:ND2	1:A:310:ASN:ND2	2.61	0.47
1:B:332:ILE:HD11	1:B:423:LEU:HD13	1.96	0.47
1:B:533:ARG:HE	1:B:534:PRO:HD2	1.80	0.47
1:B:214:ILE:O	1:B:218:VAL:HG23	2.14	0.46
1:A:521:LYS:HA	1:A:561:ARG:HG2	1.97	0.46
1:B:272:PHE:O	1:B:273:LYS:C	2.58	0.46
1:A:246:LYS:HG2	1:A:336:MET:HE3	1.97	0.46
1:A:470:LYS:HB2	1:A:470:LYS:HE3	1.70	0.46
1:B:494:GLY:O	1:B:514:ARG:NH1	2.48	0.46
1:A:277:LYS:HD2	1:A:282:LEU:HD21	1.96	0.46
1:B:433:ILE:O	1:B:437:MET:HG3	2.16	0.46
1:A:364:ASN:O	1:A:368:ARG:HG3	2.16	0.45
1:B:232:SER:OG	1:B:236:ARG:NH2	2.48	0.45
1:B:186:LEU:HD11	1:B:210:VAL:HA	1.99	0.45
1:A:345:ARG:HG3	1:A:372:TRP:CH2	2.52	0.45
1:B:246:LYS:HG2	1:B:336:MET:HE1	1.98	0.45
1:B:402:ASN:HA	1:B:405:ALA:HB3	1.99	0.45
1:A:449:LEU:HB2	1:A:489:LEU:HD11	1.99	0.44
1:A:481:GLU:HA	1:A:484:ILE:HD12	1.99	0.44
1:A:485:LYS:HG3	1:A:486:ASP:N	2.31	0.44
1:B:383:ASP:N	1:B:388:LYS:HE2	2.32	0.44
1:B:224:CYS:SG	1:B:272:PHE:O	2.76	0.44
1:A:323:PHE:CE2	1:A:324:MET:HE2	2.52	0.44
1:A:238:LEU:HD23	1:A:238:LEU:HA	1.60	0.44
1:A:265:LEU:HD23	1:A:265:LEU:HA	1.87	0.43
1:A:322:GLU:HG2	1:A:434:LEU:CD1	2.48	0.43
1:B:267:SER:OG	1:B:268:SER:N	2.49	0.43
1:A:282:LEU:HD23	1:A:282:LEU:HA	1.84	0.43
1:B:530:LEU:HD13	1:B:554:ILE:HG23	2.00	0.42
1:A:248:PRO:HD2	1:B:404:VAL:HG13	2.01	0.42
1:B:286:LYS:HG3	1:B:288:ILE:HB	2.01	0.42
1:A:187:GLU:HB2	1:A:190:ASP:OD2	2.19	0.42
1:A:348:LEU:HD13	1:A:394:LEU:HD12	2.01	0.42
1:B:226:LEU:HD12	1:B:227:ASN:H	1.84	0.42
1:B:501:PHE:HZ	1:B:554:ILE:HD13	1.84	0.42
1:B:539:ILE:HD13	1:B:539:ILE:HA	1.71	0.42
1:B:238:LEU:HA	1:B:238:LEU:HD23	1.77	0.41
1:B:435:LEU:HD13	1:B:435:LEU:HA	1.85	0.41
1:B:362:TYR:O	1:B:368:ARG:HD2	2.20	0.41
1:B:231:PHE:CE1	1:B:292:LEU:HD11	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:249:SER:HB3	1:A:250:PRO:HD2	2.02	0.41
1:B:446:LYS:HD3	1:B:448:TYR:CZ	2.55	0.41
1:B:441:LYS:HE3	1:B:441:LYS:HB2	1.91	0.41
1:A:355:GLU:HG3	1:A:356:ASN:N	2.35	0.41
1:A:393:LEU:HA	1:A:429:MET:CE	2.48	0.41
1:A:456:PRO:HD2	1:A:519:PHE:CE1	2.56	0.41
1:A:364:ASN:OD1	1:A:367:PHE:HD1	2.04	0.41
1:B:258:ILE:HG22	1:B:305:LEU:HD13	2.02	0.41
1:B:270:ARG:HG3	1:B:271:GLU:H	1.86	0.41
1:B:215:GLN:CB	1:B:231:PHE:HB2	2.51	0.40
1:A:230:LYS:HA	1:A:233:SER:OG	2.22	0.40
1:A:246:LYS:HG2	1:A:336:MET:CE	2.50	0.40
1:A:223:LEU:HD12	1:A:223:LEU:HA	1.90	0.40
1:A:259:TYR:C	1:A:259:TYR:CD1	3.00	0.40
1:A:470:LYS:HE2	1:A:557:TYR:CZ	2.57	0.40
1:B:200:ARG:HE	1:B:200:ARG:HB3	1.74	0.40
1:B:201:TYR:HB2	1:B:203:ILE:HG13	2.04	0.40
1:B:250:PRO:HD2	1:B:253:LYS:HD2	2.02	0.40
1:B:303:LEU:HA	1:B:304:PRO:HD3	1.92	0.40

All (1) 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	Interatomic distance (Å)	Clash overlap (Å)
1:A:236:ARG:NH1	1:B:500:GLU:OE1[9_555]	2.02	0.18

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	360/382 (94%)	357 (99%)	3 (1%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	356/382 (93%)	345 (97%)	11 (3%)	0	100	100
All	All	716/764 (94%)	702 (98%)	14 (2%)	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	301/360 (84%)	295 (98%)	6 (2%)	48	71
1	B	332/360 (92%)	317 (96%)	15 (4%)	24	50
All	All	633/720 (88%)	612 (97%)	21 (3%)	33	58

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	189	SER
1	A	308	LYS
1	A	318	LEU
1	A	370	LYS
1	A	510	VAL
1	A	517	ILE
1	B	217	ILE
1	B	224	CYS
1	B	232	SER
1	B	263	ASP
1	B	265	LEU
1	B	269	ILE
1	B	270	ARG
1	B	273	LYS
1	B	274	LEU
1	B	297	LYS
1	B	309	ILE
1	B	338	SER

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Mol	Chain	Res	Type
1	B	371	TYR
1	B	440	LEU
1	B	532	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	225	ASN
1	A	227	ASN
1	A	256	ASN
1	A	262	ASN
1	A	310	ASN
1	A	462	GLN
1	B	296	GLN
1	B	314	ASN
1	B	356	ASN
1	B	416	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 oligosaccharides 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 ⓘ

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	368/382 (96%)	0.19	28 (7%) 20 18	36, 55, 102, 138	0
1	B	364/382 (95%)	0.28	42 (11%) 9 9	30, 50, 120, 147	0
All	All	732/764 (95%)	0.23	70 (9%) 13 12	30, 53, 114, 147	0

All (70) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	275	CYS	8.9
1	A	533	ARG	6.4
1	B	226	LEU	6.2
1	B	379	VAL	6.0
1	B	221	ASN	6.0
1	A	376	TYR	5.3
1	B	287	PRO	5.2
1	A	185	THR	5.2
1	B	215	GLN	5.1
1	A	442	VAL	5.0
1	B	537	GLY	4.8
1	B	273	LYS	4.8
1	B	185	THR	4.7
1	B	274	LEU	4.6
1	B	378	ILE	4.4
1	B	272	PHE	4.4
1	A	227	ASN	4.1
1	A	377	ASN	4.0
1	B	271	GLU	3.9
1	A	444	SER	3.9
1	B	301	GLU	3.8
1	B	439	ASP	3.8
1	A	310	ASN	3.8
1	B	443	ASN	3.8

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Mol	Chain	Res	Type	RSRZ
1	A	481	GLU	3.7
1	A	231	PHE	3.7
1	A	539	ILE	3.6
1	B	289	GLN	3.6
1	B	219	ALA	3.6
1	B	217	ILE	3.6
1	A	543	GLU	3.6
1	B	186	LEU	3.3
1	B	187	GLU	3.3
1	B	535	GLU	3.2
1	B	229	LEU	3.1
1	B	224	CYS	3.1
1	B	376	TYR	2.9
1	A	228	GLU	2.9
1	B	230	LYS	2.9
1	B	286	LYS	2.9
1	A	367	PHE	2.9
1	A	440	LEU	2.9
1	B	228	GLU	2.7
1	B	536	GLU	2.7
1	B	225	ASN	2.7
1	A	384	LYS	2.7
1	B	383	ASP	2.6
1	B	438	LYS	2.5
1	B	213	ASN	2.5
1	B	384	LYS	2.4
1	A	550	GLU	2.4
1	A	552	ASN	2.4
1	A	229	LEU	2.4
1	B	220	LYS	2.4
1	A	439	ASP	2.3
1	A	513	LYS	2.3
1	B	231	PHE	2.3
1	A	280	GLU	2.3
1	B	216	LYS	2.2
1	A	441	LYS	2.2
1	B	392	GLU	2.2
1	A	279	SER	2.2
1	A	542	LYS	2.1
1	B	539	ILE	2.1
1	A	230	LYS	2.1
1	B	288	ILE	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	445	LYS	2.0
1	A	538	LYS	2.0
1	B	310	ASN	2.0
1	B	377	ASN	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 oligosaccharides 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.