



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 12:17 PM UTC

PDB ID : 5OJP / pdb_00005ojp
Title : YCF48 bound to D1 peptide
Authors : Michoux, F.; Nixon, P.J.; Murray, J.W.
Deposited on : 2017-07-22
Resolution : 1.86 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

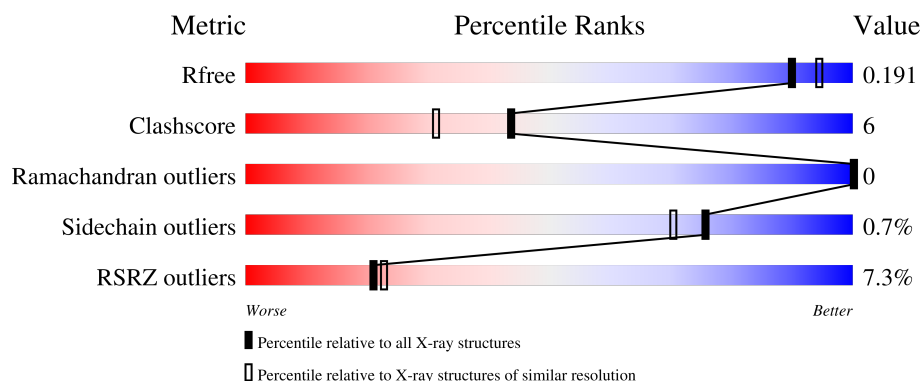
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.86 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	3428 (1.86-1.86)
Clashscore	190562	3579 (1.86-1.86)
Ramachandran outliers	187476	3553 (1.86-1.86)
Sidechain outliers	187428	3553 (1.86-1.86)
RSRZ outliers	180081	3429 (1.86-1.86)

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	326	4% 85% 10% 5%
1	B	326	2% 80% 12% 7%
1	C	326	13% 74% 18% 8%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 7331 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 Ycf48-like protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	309	Total	C	N	O	S	0	0	0
			2405	1526	409	462	8			
1	B	302	Total	C	N	O	S	0	0	0
			2360	1498	401	453	8			
1	C	299	Total	C	N	O	S	0	0	0
			2336	1485	397	446	8			

There are 63 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	22	MET	-	initiating methionine	UNP Q8DI95
A	23	ARG	-	expression tag	UNP Q8DI95
A	24	GLY	-	expression tag	UNP Q8DI95
A	25	SER	-	expression tag	UNP Q8DI95
A	26	HIS	-	expression tag	UNP Q8DI95
A	27	HIS	-	expression tag	UNP Q8DI95
A	28	HIS	-	expression tag	UNP Q8DI95
A	29	HIS	-	expression tag	UNP Q8DI95
A	30	HIS	-	expression tag	UNP Q8DI95
A	31	HIS	-	expression tag	UNP Q8DI95
A	32	GLY	-	expression tag	UNP Q8DI95
A	33	LEU	-	expression tag	UNP Q8DI95
A	34	VAL	-	expression tag	UNP Q8DI95
A	35	PRO	-	expression tag	UNP Q8DI95
A	36	ARG	-	expression tag	UNP Q8DI95
A	37	GLY	-	expression tag	UNP Q8DI95
A	38	SER	-	expression tag	UNP Q8DI95
A	202	GLU	ARG	engineered mutation	UNP Q8DI95
A	221	GLU	ARG	engineered mutation	UNP Q8DI95
A	225	GLU	ARG	engineered mutation	UNP Q8DI95
A	226	GLU	ARG	engineered mutation	UNP Q8DI95
B	22	MET	-	initiating methionine	UNP Q8DI95
B	23	ARG	-	expression tag	UNP Q8DI95

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Chain	Residue	Modelled	Actual	Comment	Reference
B	24	GLY	-	expression tag	UNP Q8DI95
B	25	SER	-	expression tag	UNP Q8DI95
B	26	HIS	-	expression tag	UNP Q8DI95
B	27	HIS	-	expression tag	UNP Q8DI95
B	28	HIS	-	expression tag	UNP Q8DI95
B	29	HIS	-	expression tag	UNP Q8DI95
B	30	HIS	-	expression tag	UNP Q8DI95
B	31	HIS	-	expression tag	UNP Q8DI95
B	32	GLY	-	expression tag	UNP Q8DI95
B	33	LEU	-	expression tag	UNP Q8DI95
B	34	VAL	-	expression tag	UNP Q8DI95
B	35	PRO	-	expression tag	UNP Q8DI95
B	36	ARG	-	expression tag	UNP Q8DI95
B	37	GLY	-	expression tag	UNP Q8DI95
B	38	SER	-	expression tag	UNP Q8DI95
B	202	GLU	ARG	engineered mutation	UNP Q8DI95
B	221	GLU	ARG	engineered mutation	UNP Q8DI95
B	225	GLU	ARG	engineered mutation	UNP Q8DI95
B	226	GLU	ARG	engineered mutation	UNP Q8DI95
C	22	MET	-	initiating methionine	UNP Q8DI95
C	23	ARG	-	expression tag	UNP Q8DI95
C	24	GLY	-	expression tag	UNP Q8DI95
C	25	SER	-	expression tag	UNP Q8DI95
C	26	HIS	-	expression tag	UNP Q8DI95
C	27	HIS	-	expression tag	UNP Q8DI95
C	28	HIS	-	expression tag	UNP Q8DI95
C	29	HIS	-	expression tag	UNP Q8DI95
C	30	HIS	-	expression tag	UNP Q8DI95
C	31	HIS	-	expression tag	UNP Q8DI95
C	32	GLY	-	expression tag	UNP Q8DI95
C	33	LEU	-	expression tag	UNP Q8DI95
C	34	VAL	-	expression tag	UNP Q8DI95
C	35	PRO	-	expression tag	UNP Q8DI95
C	36	ARG	-	expression tag	UNP Q8DI95
C	37	GLY	-	expression tag	UNP Q8DI95
C	38	SER	-	expression tag	UNP Q8DI95
C	202	GLU	ARG	engineered mutation	UNP Q8DI95
C	221	GLU	ARG	engineered mutation	UNP Q8DI95
C	225	GLU	ARG	engineered mutation	UNP Q8DI95
C	226	GLU	ARG	engineered mutation	UNP Q8DI95

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	97	Total 97	O 97	0	0
2	B	91	Total 91	O 91	0	0
2	C	42	Total 42	O 42	0	0

4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	54.44Å 60.29Å 75.40Å 68.43° 74.83° 85.43°	Depositor
Resolution (Å)	53.77 – 1.86 53.77 – 1.86	Depositor EDS
% Data completeness (in resolution range)	82.8 (53.77-1.86) 83.0 (53.77-1.86)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.52 (at 1.86Å)	Xtriage
Refinement program	PHENIX (1.12_2829: ???)	Depositor
R, R_{free}	0.190 , 0.236 0.194 , 0.191	Depositor DCC
R_{free} test set	3016 reflections (4.17%)	wwPDB-VP
Wilson B-factor (Å ²)	31.9	Xtriage
Anisotropy	0.272	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 43.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	7331	wwPDB-VP
Average B, all atoms (Å ²)	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.86% 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	0.68	0/2473	0.85	0/3375
1	B	0.69	0/2426	0.84	0/3307
1	C	0.56	0/2401	0.76	0/3271
All	All	0.65	0/7300	0.82	0/9953

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	B	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	338	TYR	Sidechain

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	2405	0	2297	21	1
1	B	2360	0	2253	26	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	2336	0	2230	39	1
2	A	97	0	0	0	0
2	B	91	0	0	4	0
2	C	42	0	0	0	0
All	All	7331	0	6780	84	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 6.

All (84) 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:C:222:THR:HG21	1:C:240:MET:HE1	1.32	1.12
1:A:58:ILE:O	1:A:316:TYR:OH	1.72	1.06
1:B:317:LYS:HE2	1:B:319:LEU:HD21	1.65	0.77
1:B:314:ASN:HB2	1:B:316:TYR:CE1	2.27	0.70
1:C:40:PRO:HG2	1:C:303:GLN:HG2	1.74	0.70
1:C:237:ARG:HD2	1:C:250:SER:O	1.94	0.68
1:C:150:LEU:HD11	1:C:156:GLU:HG3	1.77	0.67
1:B:38:SER:N	2:B:402:HOH:O	2.27	0.67
1:A:317:LYS:HE2	1:A:319:LEU:HD21	1.77	0.66
1:C:304:GLN:NE2	1:C:309:LYS:HD2	2.11	0.65
1:C:184:VAL:O	1:C:201:SER:HB3	2.03	0.59
1:B:300:GLN:NE2	2:B:404:HOH:O	2.36	0.58
1:B:226:GLU:HB3	1:B:243:ASN:HB2	1.86	0.58
1:A:158:ILE:HG12	1:A:164:ILE:HG12	1.86	0.58
1:C:117:GLU:HA	1:C:118:PRO:C	2.29	0.58
1:A:345:ALA:HA	1:C:273:LEU:HD21	1.86	0.57
1:C:158:ILE:HG12	1:C:164:ILE:HG12	1.86	0.57
1:A:328:ILE:HB	1:A:336:LEU:HB2	1.87	0.56
1:C:251:ASP:HB2	1:C:257:ASN:HB2	1.86	0.56
1:C:157:MET:HG2	1:C:158:ILE:N	2.22	0.55
1:A:150:LEU:HD11	1:A:156:GLU:HG3	1.89	0.55
1:C:163:ALA:HB1	1:C:165:TYR:HE1	1.71	0.55
1:B:150:LEU:HD11	1:B:156:GLU:HG3	1.89	0.54
1:B:117:GLU:HA	1:B:118:PRO:C	2.32	0.54
1:C:318:ILE:HG12	1:C:328:ILE:HG12	1.90	0.54
1:C:144:PRO:HB3	1:C:157:MET:SD	2.49	0.53
1:B:71:LEU:C	1:B:71:LEU:HD12	2.34	0.53
1:C:286:ALA:HB2	1:C:318:ILE:HD11	1.90	0.53
1:A:163:ALA:HB1	1:A:165:TYR:HE1	1.74	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:101:ARG:HD3	1:A:117:GLU:OE2	2.08	0.52
1:C:240:MET:HE3	1:C:258:TRP:CZ2	2.45	0.52
1:B:227:LEU:HD23	1:B:242:VAL:HG12	1.92	0.52
1:B:314:ASN:HB2	1:B:316:TYR:HE1	1.73	0.50
1:C:164:ILE:C	1:C:165:TYR:HD1	2.19	0.50
1:C:233:THR:OG1	1:C:235:ASP:OD1	2.26	0.50
1:C:48:GLU:HG3	1:C:339:VAL:HG21	1.94	0.50
1:B:337:ARG:HD2	2:B:422:HOH:O	2.11	0.50
1:B:86:GLN:HG2	1:B:337:ARG:CZ	2.42	0.49
1:C:115:VAL:HB	1:C:144:PRO:HB2	1.93	0.49
1:A:117:GLU:HA	1:A:118:PRO:C	2.38	0.48
1:C:328:ILE:HB	1:C:336:LEU:HB2	1.95	0.48
1:B:55:THR:O	1:B:332:LYS:HA	2.13	0.48
1:B:184:VAL:O	1:B:201:SER:HB3	2.14	0.48
1:C:157:MET:HE2	1:C:157:MET:HB3	1.75	0.47
1:A:163:ALA:HB1	1:A:165:TYR:CE1	2.50	0.47
1:B:275:LEU:C	1:B:275:LEU:HD23	2.40	0.46
1:C:39:ILE:HB	1:C:40:PRO:HD2	1.97	0.46
1:B:265:LEU:HG	1:B:293:LEU:HD11	1.97	0.46
1:B:324:ASP:N	1:B:324:ASP:OD1	2.48	0.46
1:C:246:LYS:HA	1:C:262:LEU:O	2.16	0.45
1:C:317:LYS:HE2	1:C:319:LEU:HD21	1.97	0.45
1:A:240:MET:HE3	1:A:258:TRP:CD2	2.51	0.45
1:A:137:ASP:OD1	1:A:138:PRO:HD2	2.17	0.45
1:A:96:ASP:N	1:A:96:ASP:OD1	2.48	0.44
1:B:185:MET:HE3	1:B:185:MET:HB2	1.81	0.44
1:B:237:ARG:NH1	2:B:401:HOH:O	2.16	0.44
1:B:58:ILE:HD12	1:B:71:LEU:HD13	2.00	0.44
1:C:237:ARG:HG2	1:C:252:PRO:HA	2.00	0.43
1:C:290:GLY:HA2	1:C:315:PHE:CE2	2.53	0.43
1:C:122:LEU:HD23	1:C:133:GLN:HA	1.99	0.43
1:C:45:ASN:HB2	1:C:307:ASP:OD2	2.18	0.43
1:C:304:GLN:HE22	1:C:309:LYS:HD2	1.83	0.43
1:B:297:ASP:OD2	1:B:300:GLN:HB2	2.19	0.42
1:A:115:VAL:HB	1:A:144:PRO:HB2	2.01	0.42
1:B:328:ILE:HB	1:B:336:LEU:HB2	2.02	0.42
1:A:58:ILE:HD12	1:A:71:LEU:HD13	2.01	0.42
1:C:58:ILE:HD12	1:C:71:LEU:HD13	2.02	0.42
1:C:221:GLU:HG2	1:C:222:THR:HG23	2.02	0.42
1:A:211:PRO:HB2	1:B:211:PRO:HB2	2.02	0.42
1:C:146:LEU:HD23	1:C:147:ILE:N	2.35	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:186:ARG:NH1	1:C:227:LEU:O	2.53	0.41
1:C:207:SER:HA	1:C:217:GLU:O	2.20	0.41
1:C:146:LEU:HD23	1:C:146:LEU:C	2.45	0.41
1:A:146:LEU:C	1:A:146:LEU:HD23	2.45	0.41
1:A:71:LEU:HD12	1:A:71:LEU:C	2.45	0.41
1:B:115:VAL:HB	1:B:144:PRO:HB2	2.03	0.41
1:C:190:ARG:CZ	1:C:194:GLY:HA2	2.51	0.41
1:C:79:MET:HG2	1:C:90:PRO:HA	2.03	0.41
1:A:282:GLU:HA	1:A:295:SER:O	2.21	0.40
1:A:327:PHE:HA	1:A:336:LEU:O	2.21	0.40
1:B:277:TYR:CE2	1:B:283:VAL:HG22	2.57	0.40
1:C:221:GLU:HG2	1:C:222:THR:N	2.35	0.40
1:A:58:ILE:O	1:A:316:TYR:CZ	2.69	0.40
1:B:39:ILE:HB	1:B:40:PRO:HD2	2.03	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:220:ASN:O	1:C:172:LYS:NZ[1_564]	2.04	0.16

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	307/326 (94%)	303 (99%)	4 (1%)	0	100	100
1	B	298/326 (91%)	297 (100%)	1 (0%)	0	100	100
1	C	293/326 (90%)	291 (99%)	2 (1%)	0	100	100
All	All	898/978 (92%)	891 (99%)	7 (1%)	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	258/272 (95%)	257 (100%)	1 (0%)	84	82
1	B	254/272 (93%)	251 (99%)	3 (1%)	63	55
1	C	250/272 (92%)	249 (100%)	1 (0%)	84	82
All	All	762/816 (93%)	757 (99%)	5 (1%)	76	70

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	324	ASP
1	B	165	TYR
1	B	316	TYR
1	B	338	TYR
1	C	165	TYR

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

Mol	Chain	Res	Type
1	A	108	GLN
1	A	129	GLN
1	A	220	ASN
1	B	129	GLN
1	B	213	GLN
1	B	257	ASN
1	B	331	GLN
1	C	110	ASN

5.3.3 RNA ⓘ

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 [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	309/326 (94%)	0.38	14 (4%) 38 41	23, 40, 73, 104	0
1	B	302/326 (92%)	0.35	8 (2%) 57 61	23, 42, 77, 119	0
1	C	299/326 (91%)	1.04	44 (14%) 6 5	31, 57, 111, 143	0
All	All	910/978 (93%)	0.59	66 (7%) 21 23	23, 46, 91, 143	0

All (66) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	316	TYR	5.4
1	C	223	THR	4.4
1	C	222	THR	4.3
1	C	212	GLY	4.3
1	A	346	PRO	3.7
1	C	267	ARG	3.6
1	C	252	PRO	3.5
1	C	141	PRO	3.2
1	A	315	PHE	3.2
1	B	269	SER	3.1
1	C	249	PHE	3.1
1	B	44	TYR	3.1
1	A	342	LEU	3.0
1	C	199	VAL	3.0
1	C	272	PHE	3.0
1	C	178	VAL	3.0
1	C	218	PRO	3.0
1	C	142	GLY	2.9
1	C	316	TYR	2.9
1	C	228	HIS	2.9
1	A	344	ALA	2.9
1	C	140	LEU	2.9
1	C	261	LEU	2.9

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Mol	Chain	Res	Type	RSRZ
1	C	265	LEU	2.9
1	C	271	GLY	2.9
1	C	184	VAL	2.8
1	C	255	SER	2.8
1	C	312	PRO	2.8
1	C	101	ARG	2.7
1	B	266	ARG	2.6
1	C	238	LEU	2.6
1	C	289	ALA	2.6
1	B	340	THR	2.6
1	C	44	TYR	2.5
1	C	136	LEU	2.5
1	C	214	THR	2.5
1	C	236	GLY	2.4
1	A	180	GLU	2.4
1	A	313	SER	2.4
1	B	38	SER	2.4
1	A	179	GLN	2.3
1	C	260	GLU	2.3
1	C	227	LEU	2.3
1	C	333	GLY	2.3
1	B	92	THR	2.3
1	A	312	PRO	2.3
1	C	135	PRO	2.3
1	C	245	GLY	2.3
1	C	243	ASN	2.3
1	C	266	ARG	2.2
1	C	215	ALA	2.2
1	C	291	ALA	2.2
1	C	216	TRP	2.2
1	C	182	ILE	2.2
1	A	314	ASN	2.2
1	B	341	ASP	2.1
1	C	54	THR	2.1
1	C	134	ILE	2.1
1	A	343	THR	2.1
1	A	94	VAL	2.1
1	A	345	ALA	2.1
1	C	220	ASN	2.0
1	C	41	ALA	2.0
1	A	260	GLU	2.0
1	B	39	ILE	2.0

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Mol	Chain	Res	Type	RSRZ
1	C	165	TYR	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.