



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 8, 2021 – 02:17 PM JST

PDB ID : 7BUW
Title : Eucommia ulmoides TPT3 mutant -C94Y/A95F
Authors : Kajiura, H.; Yoshizawa, T.; Tokumoto, Y.; Suzuki, N.; Takeno, S.; Takeno, K.J.; Yamashita, T.; Tanaka, S.; Kaneko, Y.; Fujiyama, K.; Matsumura, H.; Nakazawa, Y.
Deposited on : 2020-04-08
Resolution : 3.30 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	:	2.16
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)	:	2.16

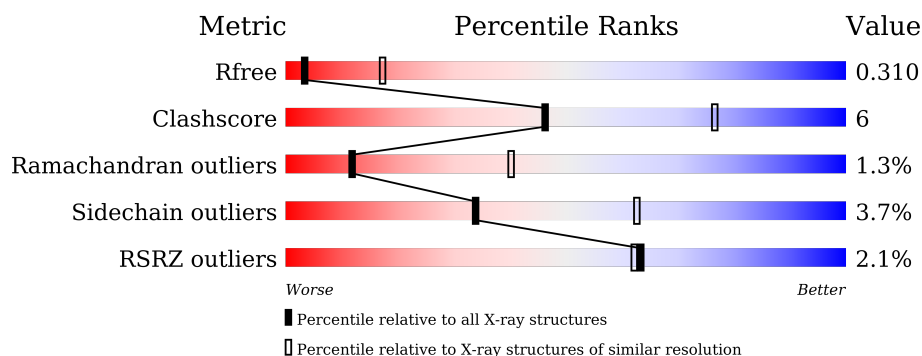
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.30 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	1149 (3.34-3.26)
Clashscore	141614	1205 (3.34-3.26)
Ramachandran outliers	138981	1183 (3.34-3.26)
Sidechain outliers	138945	1182 (3.34-3.26)
RSRZ outliers	127900	1115 (3.34-3.26)

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	362	 72% 13% • 13%
1	B	362	 73% 12% • 13%

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 5108 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 FPS3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	314	Total	C	N	O	S	0	0	0
			2550	1640	412	484	14			
1	B	315	Total	C	N	O	S	0	0	0
			2558	1644	415	486	13			

There are 44 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-15	MET	-	initiating methionine	UNP A0A1L3KPU1
A	-14	ASN	-	expression tag	UNP A0A1L3KPU1
A	-13	HIS	-	expression tag	UNP A0A1L3KPU1
A	-12	LYS	-	expression tag	UNP A0A1L3KPU1
A	-11	VAL	-	expression tag	UNP A0A1L3KPU1
A	-10	HIS	-	expression tag	UNP A0A1L3KPU1
A	-9	HIS	-	expression tag	UNP A0A1L3KPU1
A	-8	HIS	-	expression tag	UNP A0A1L3KPU1
A	-7	HIS	-	expression tag	UNP A0A1L3KPU1
A	-6	HIS	-	expression tag	UNP A0A1L3KPU1
A	-5	HIS	-	expression tag	UNP A0A1L3KPU1
A	-4	ILE	-	expression tag	UNP A0A1L3KPU1
A	-3	GLU	-	expression tag	UNP A0A1L3KPU1
A	-2	GLY	-	expression tag	UNP A0A1L3KPU1
A	-1	ARG	-	expression tag	UNP A0A1L3KPU1
A	0	HIS	-	expression tag	UNP A0A1L3KPU1
A	68	GLY	ASN	engineered mutation	UNP A0A1L3KPU1
A	69	GLY	SER	engineered mutation	UNP A0A1L3KPU1
A	?	-	LYS	deletion	UNP A0A1L3KPU1
A	?	-	TYR	deletion	UNP A0A1L3KPU1
A	94	TYR	CYS	engineered mutation	UNP A0A1L3KPU1
A	95	PHE	ALA	engineered mutation	UNP A0A1L3KPU1
B	-15	MET	-	initiating methionine	UNP A0A1L3KPU1
B	-14	ASN	-	expression tag	UNP A0A1L3KPU1
B	-13	HIS	-	expression tag	UNP A0A1L3KPU1

Continued on next page...

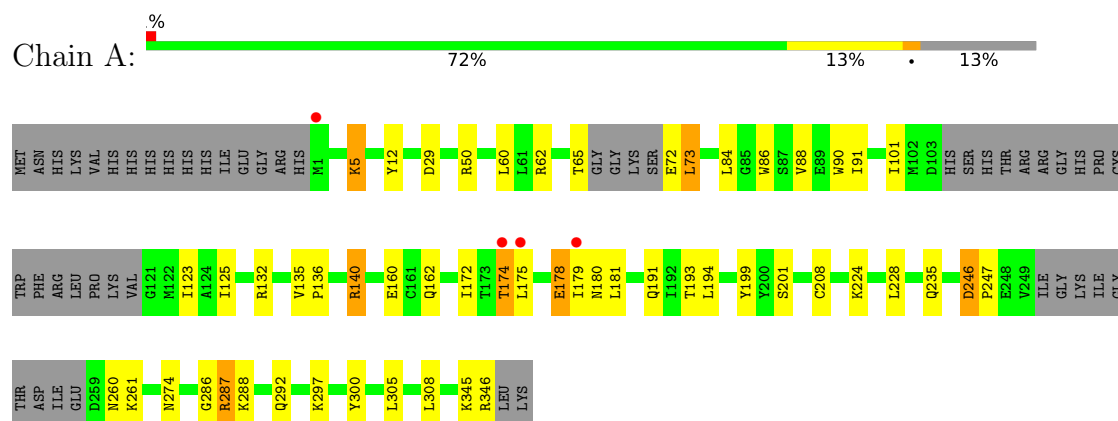
Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
B	-12	LYS	-	expression tag	UNP A0A1L3KPU1
B	-11	VAL	-	expression tag	UNP A0A1L3KPU1
B	-10	HIS	-	expression tag	UNP A0A1L3KPU1
B	-9	HIS	-	expression tag	UNP A0A1L3KPU1
B	-8	HIS	-	expression tag	UNP A0A1L3KPU1
B	-7	HIS	-	expression tag	UNP A0A1L3KPU1
B	-6	HIS	-	expression tag	UNP A0A1L3KPU1
B	-5	HIS	-	expression tag	UNP A0A1L3KPU1
B	-4	ILE	-	expression tag	UNP A0A1L3KPU1
B	-3	GLU	-	expression tag	UNP A0A1L3KPU1
B	-2	GLY	-	expression tag	UNP A0A1L3KPU1
B	-1	ARG	-	expression tag	UNP A0A1L3KPU1
B	0	HIS	-	expression tag	UNP A0A1L3KPU1
B	68	GLY	ASN	engineered mutation	UNP A0A1L3KPU1
B	69	GLY	SER	engineered mutation	UNP A0A1L3KPU1
B	?	-	LYS	deletion	UNP A0A1L3KPU1
B	?	-	TYR	deletion	UNP A0A1L3KPU1
B	94	TYR	CYS	engineered mutation	UNP A0A1L3KPU1
B	95	PHE	ALA	engineered mutation	UNP A0A1L3KPU1

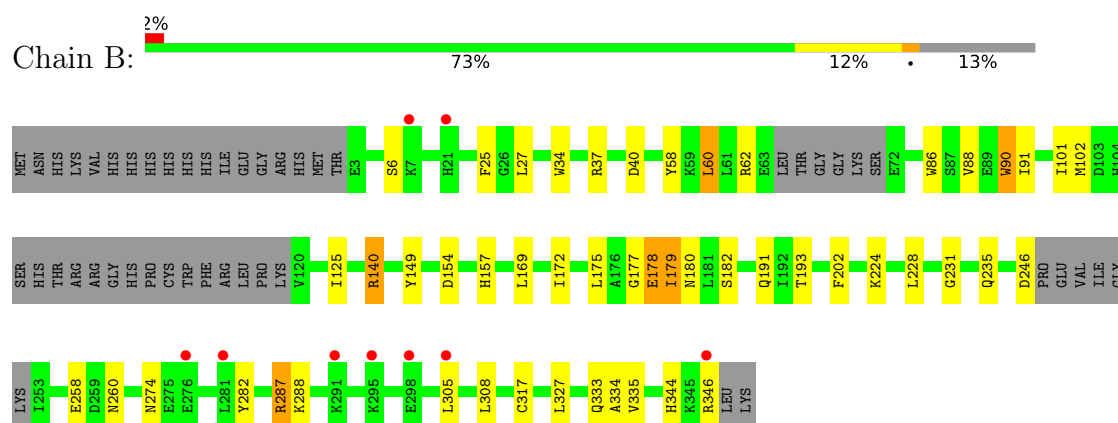
3 Residue-property plots [i](#)

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: FPS3



• Molecule 1: FPS3



4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	84.25Å 84.25Å 181.19Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	46.53 – 3.30 46.53 – 3.20	Depositor EDS
% Data completeness (in resolution range)	100.0 (46.53-3.30) 100.0 (46.53-3.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.06 (at 3.19Å)	Xtriage
Refinement program	PHENIX 1.17.1_3660	Depositor
R, R_{free}	0.253 , 0.310 0.253 , 0.310	Depositor DCC
R_{free} test set	588 reflections (4.57%)	wwPDB-VP
Wilson B-factor (Å ²)	68.2	Xtriage
Anisotropy	0.440	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 32.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.039 for -h,-k,l	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	5108	wwPDB-VP
Average B, all atoms (Å ²)	65.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.88% 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	0.29	0/2597	0.48	0/3511
1	B	0.28	0/2605	0.46	0/3521
All	All	0.29	0/5202	0.47	0/7032

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

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	2550	0	2552	32	0
1	B	2558	0	2551	38	0
All	All	5108	0	5103	64	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 6.

All (64) 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:60:LEU:HD22	1:B:335:VAL:HG22	1.28	1.15
1:B:60:LEU:HD22	1:B:335:VAL:CG2	1.76	1.15

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:60:LEU:CD2	1:B:335:VAL:HG23	1.97	0.95
1:B:60:LEU:CD2	1:B:335:VAL:CG2	2.45	0.93
1:B:287:ARG:HH11	1:B:287:ARG:HG2	1.38	0.87
1:B:178:GLU:HG2	1:B:179:ILE:HG13	1.59	0.82
1:A:345:LYS:O	1:A:346:ARG:HG2	1.80	0.82
1:B:287:ARG:NH1	1:B:287:ARG:HG2	1.98	0.74
1:B:60:LEU:HD23	1:B:335:VAL:HG23	1.68	0.73
1:A:136:PRO:HB2	1:B:157:HIS:CE1	2.29	0.67
1:B:60:LEU:CD2	1:B:335:VAL:HG22	2.12	0.65
1:B:179:ILE:HD12	1:B:180:ASN:N	2.13	0.64
1:B:101:ILE:HD13	1:B:125:ILE:HD13	1.79	0.64
1:A:123:ILE:HD11	1:B:175:LEU:HD11	1.83	0.61
1:B:193:THR:HG22	1:B:235:GLN:HG2	1.81	0.60
1:A:72:GLU:O	1:A:73:LEU:HB2	2.01	0.60
1:A:178:GLU:HB3	1:A:179:ILE:HG13	1.85	0.59
1:A:193:THR:HG22	1:A:235:GLN:HG2	1.84	0.59
1:B:175:LEU:O	1:B:178:GLU:HB3	2.02	0.58
1:A:140:ARG:HH11	1:B:154:ASP:CG	2.07	0.57
1:A:345:LYS:O	1:A:346:ARG:CG	2.54	0.55
1:A:101:ILE:HD13	1:A:125:ILE:HD13	1.89	0.55
1:B:178:GLU:CG	1:B:179:ILE:HG13	2.35	0.54
1:B:37:ARG:NH2	1:B:40:ASP:OD2	2.40	0.54
1:A:174:THR:HG22	1:A:175:LEU:N	2.22	0.53
1:B:177:GLY:O	1:B:260:ASN:ND2	2.40	0.51
1:A:140:ARG:NH1	1:B:154:ASP:CG	2.64	0.50
1:B:287:ARG:CG	1:B:287:ARG:HH11	2.14	0.50
1:B:327:LEU:O	1:B:333:GLN:NE2	2.42	0.50
1:A:180:ASN:OD1	1:A:180:ASN:N	2.45	0.48
1:B:60:LEU:HD21	1:B:334:ALA:HB3	1.96	0.48
1:A:305:LEU:HA	1:A:308:LEU:HB2	1.95	0.47
1:A:287:ARG:HD2	1:A:292:GLN:HE22	1.78	0.47
1:A:287:ARG:HD2	1:A:292:GLN:NE2	2.29	0.47
1:A:88:VAL:O	1:A:91:ILE:HG22	2.15	0.46
1:B:258:GLU:HB3	1:B:288:LYS:HG2	1.97	0.46
1:A:84:LEU:HD13	1:A:208:CYS:HB3	1.98	0.46
1:B:88:VAL:O	1:B:91:ILE:HG22	2.16	0.45
1:A:140:ARG:NH1	1:B:154:ASP:HB3	2.31	0.45
1:A:288:LYS:HB3	1:A:288:LYS:HE3	1.83	0.45
1:A:91:ILE:HD12	1:A:91:ILE:HA	1.81	0.45
1:A:62:ARG:HA	1:A:65:THR:HG22	1.99	0.45
1:A:172:ILE:HD11	1:B:34:TRP:CG	2.53	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:12:TYR:HB2	1:A:86:TRP:CZ2	2.53	0.44
1:A:224:LYS:O	1:A:228:LEU:HD13	2.18	0.43
1:A:5:LYS:HD2	1:A:5:LYS:C	2.38	0.43
1:A:132:ARG:NH1	1:A:160:GLU:OE1	2.51	0.43
1:B:140:ARG:HG3	1:B:149:TYR:OH	2.19	0.43
1:A:162:GLN:OE1	1:A:199:TYR:CZ	2.72	0.42
1:B:169:LEU:HD23	1:B:172:ILE:HD11	2.01	0.42
1:B:224:LYS:O	1:B:228:LEU:HD13	2.19	0.42
1:B:317:CYS:SG	1:B:344:HIS:ND1	2.92	0.42
1:B:86:TRP:O	1:B:90:TRP:HE3	2.02	0.42
1:A:297:LYS:O	1:A:300:TYR:HB2	2.19	0.42
1:B:58:TYR:O	1:B:62:ARG:HG2	2.20	0.42
1:B:60:LEU:HD21	1:B:334:ALA:CB	2.49	0.41
1:B:305:LEU:HA	1:B:308:LEU:HB2	2.03	0.41
1:B:180:ASN:OD1	1:B:182:SER:HB2	2.19	0.41
1:B:202:PHE:CE1	1:B:231:GLY:HA2	2.55	0.41
1:A:50:ARG:HE	1:A:201:SER:HA	1.85	0.41
1:A:260:ASN:HB2	1:A:286:GLY:HA2	2.03	0.41
1:B:25:PHE:O	1:B:27:LEU:HD12	2.21	0.41
1:A:135:VAL:HB	1:A:136:PRO:HD3	2.02	0.41
1:A:194:LEU:CD2	1:A:228:LEU:HB3	2.52	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	306/362 (84%)	281 (92%)	19 (6%)	6 (2%)	7	32
1	B	307/362 (85%)	288 (94%)	17 (6%)	2 (1%)	22	54
All	All	613/724 (85%)	569 (93%)	36 (6%)	8 (1%)	12	40

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	174	THR
1	A	178	GLU
1	A	246	ASP
1	A	247	PRO
1	A	73	LEU
1	B	274	ASN
1	A	274	ASN
1	B	179	ILE

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	287/329 (87%)	277 (96%)	10 (4%)	36	64
1	B	287/329 (87%)	276 (96%)	11 (4%)	33	62
All	All	574/658 (87%)	553 (96%)	21 (4%)	34	63

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

Mol	Chain	Res	Type
1	A	5	LYS
1	A	29	ASP
1	A	60	LEU
1	A	90	TRP
1	A	140	ARG
1	A	181	LEU
1	A	191	GLN
1	A	246	ASP
1	A	261	LYS
1	A	287	ARG
1	B	6	SER
1	B	60	LEU
1	B	90	TRP
1	B	102	MET
1	B	140	ARG

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	178	GLU
1	B	191	GLN
1	B	246	ASP
1	B	282	TYR
1	B	287	ARG
1	B	346	ARG

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

Mol	Chain	Res	Type
1	B	157	HIS

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, 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	314/362 (86%)	-0.17	4 (1%) 77 77	33, 52, 84, 117	0
1	B	315/362 (87%)	0.13	9 (2%) 51 50	43, 72, 106, 128	0
All	All	629/724 (86%)	-0.02	13 (2%) 63 62	33, 62, 98, 128	0

All (13) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	175	LEU	5.4
1	A	179	ILE	3.3
1	B	298	GLU	2.9
1	A	174	THR	2.8
1	B	295	LYS	2.6
1	A	1	MET	2.5
1	B	291	LYS	2.3
1	B	21	HIS	2.3
1	B	276	GLU	2.3
1	B	305	LEU	2.3
1	B	346	ARG	2.1
1	B	281	LEU	2.0
1	B	7	LYS	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.