Full wwPDB X-ray Structure Validation Report (i)

Feb 19, 2016 – 10:52 PM GMT

PDB ID	:	5ERA
Title	:	Human Connexin-26 (Calcium-free)
Authors	:	Purdy, M.D.; Bennett, B.C.; Baker, K.A.; Yeager, M.J.
Deposited on	:	2015-11-13
Resolution	:	3.80 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 (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

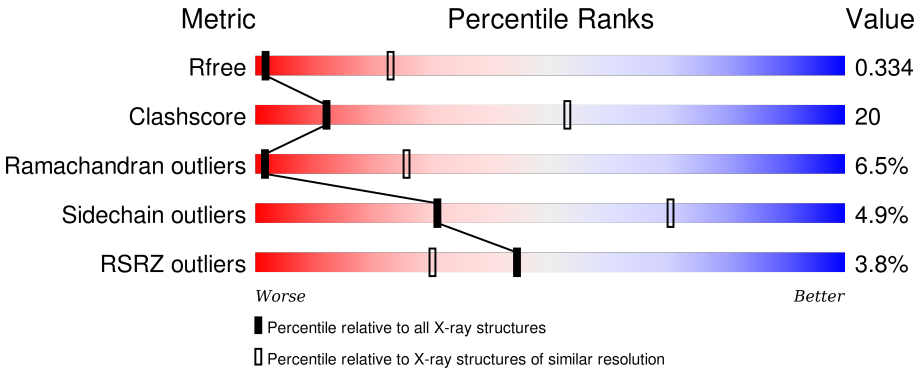
MolProbity	:	4.02b-467
Mogul	:	unknown
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026982
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	rb-20026982

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 3.80 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	1317 (4.10-3.50)
Clashscore	102246	1458 (4.10-3.50)
Ramachandran outliers	100387	1397 (4.10-3.50)
Sidechain outliers	100360	1392 (4.10-3.50)
RSRZ outliers	91569	1325 (4.10-3.50)

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 <=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	226	% 54% 11% 29%
1	B	226	4% 55% 9% 31%

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 2539 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 Gap junction beta-2 protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	155	Total	C	N	O	S	0	0	0
			1251	835	194	209	13			
1	A	160	Total	C	N	O	S	0	0	0
			1288	860	199	216	13			

There are 6 discrepancies between the modelled and reference sequences:

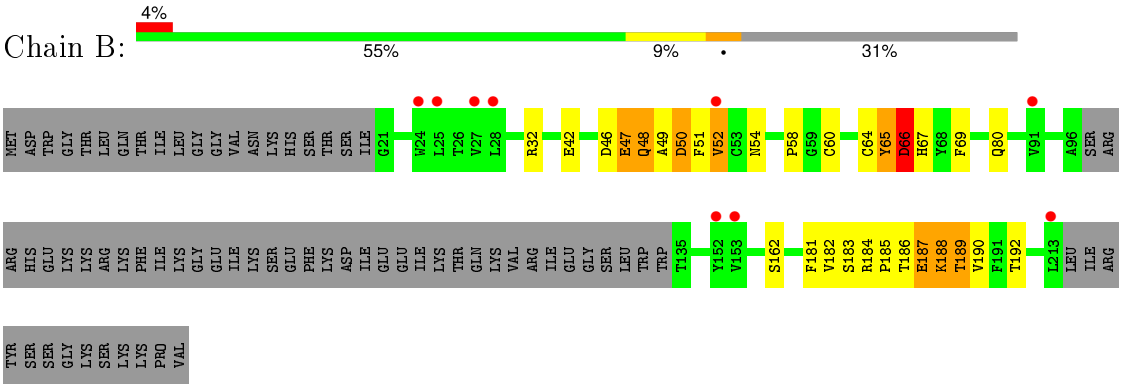
Chain	Residue	Modelled	Actual	Comment	Reference
B	97	SER	TYR	engineered mutation	UNP P29033
B	211	SER	CYS	engineered mutation	UNP P29033
B	218	SER	CYS	engineered mutation	UNP P29033
A	97	SER	TYR	engineered mutation	UNP P29033
A	211	SER	CYS	engineered mutation	UNP P29033
A	218	SER	CYS	engineered mutation	UNP P29033

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.

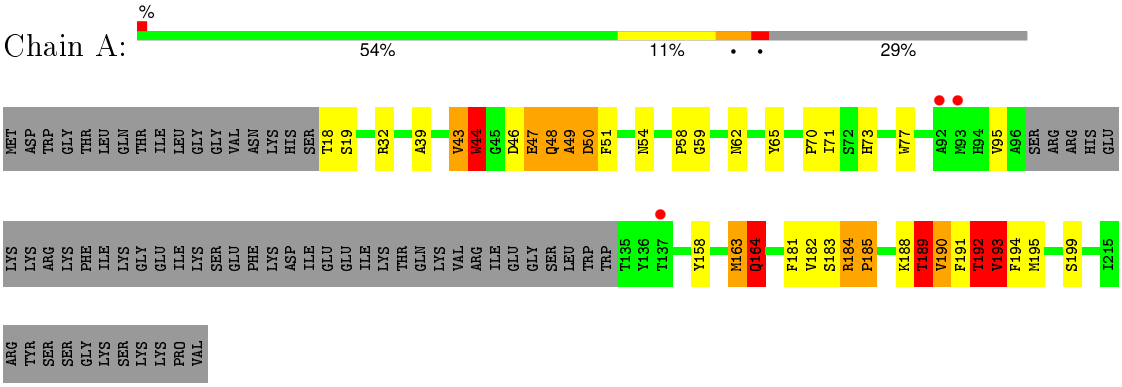
- Molecule 1: Gap junction beta-2 protein

Chain B:



- Molecule 1: Gap junction beta-2 protein

Chain A:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants a, b, c, α, β, γ	155.73Å 155.73Å 160.70Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	44.13 – 3.80 44.13 – 3.51	Depositor EDS
% Data completeness (in resolution range)	97.9 (44.13-3.80) 85.9 (44.13-3.51)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.42 (at 3.48Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.300 , 0.333 0.302 , 0.334	Depositor DCC
R_{free} test set	685 reflections (9.99%)	DCC
Wilson B-factor (Å^2)	125.1	Xtriage
Anisotropy	0.223	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.29 , 96.8	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.47$, $< L^2 > = 0.30$	Xtriage
Outliers	3 of 9419 reflections (0.032%)	Xtriage
F_o,F_c correlation	0.84	EDS
Total number of atoms	2539	wwPDB-VP
Average B, all atoms (Å^2)	127.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.64% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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	0.38	0/1324	0.58	0/1805
1	B	0.34	0/1287	0.56	0/1754
All	All	0.36	0/2611	0.57	0/3559

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	1288	0	1297	71	1
1	B	1251	0	1252	40	1
All	All	2539	0	2549	101	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 20.

All (101) 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:181:PHE:HE1	1:A:58:PRO:O	1.13	1.25
1:B:181:PHE:CE1	1:A:58:PRO:O	1.98	1.14

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:193:VAL:HG12	1:A:194:PHE:N	1.70	1.04
1:A:193:VAL:CG1	1:A:194:PHE:N	2.32	0.91
1:A:44:TRP:HZ2	1:A:73:HIS:CG	1.90	0.89
1:A:193:VAL:HG12	1:A:194:PHE:H	1.40	0.87
1:A:47:GLU:O	1:A:50:ASP:N	2.07	0.86
1:A:51:PHE:HE2	1:A:65:TYR:HB2	1.40	0.86
1:B:181:PHE:CD1	1:A:59:GLY:HA3	2.11	0.85
1:B:65:TYR:CE2	1:B:69:PHE:HB2	2.14	0.83
1:A:51:PHE:CE2	1:A:65:TYR:HB2	2.13	0.83
1:A:44:TRP:CZ2	1:A:73:HIS:HB2	2.15	0.81
1:A:190:VAL:HG12	1:A:191:PHE:N	1.95	0.80
1:A:39:ALA:HB2	1:A:191:PHE:HD1	1.48	0.77
1:B:65:TYR:CD2	1:B:69:PHE:HB2	2.21	0.76
1:A:77:TRP:HZ2	1:A:192:THR:HG22	1.49	0.76
1:A:44:TRP:CZ2	1:A:73:HIS:CG	2.73	0.76
1:A:48:GLN:O	1:A:49:ALA:C	2.24	0.72
1:B:65:TYR:CE2	1:B:69:PHE:CB	2.73	0.72
1:A:190:VAL:O	1:A:191:PHE:C	2.29	0.71
1:B:181:PHE:CE1	1:A:59:GLY:HA3	2.25	0.70
1:A:48:GLN:O	1:A:50:ASP:N	2.25	0.70
1:A:43:VAL:HG12	1:A:44:TRP:N	2.06	0.70
1:A:43:VAL:O	1:A:44:TRP:C	2.30	0.69
1:A:192:THR:O	1:A:193:VAL:C	2.29	0.69
1:A:192:THR:O	1:A:195:MET:N	2.25	0.69
1:A:77:TRP:CZ2	1:A:192:THR:HG22	2.26	0.69
1:A:44:TRP:CZ2	1:A:73:HIS:CB	2.76	0.69
1:A:47:GLU:O	1:A:48:GLN:C	2.29	0.69
1:A:188:LYS:O	1:A:189:THR:C	2.30	0.69
1:A:48:GLN:O	1:A:51:PHE:N	2.27	0.67
1:B:32:ARG:NH2	1:B:80:GLN:OE1	2.28	0.66
1:A:46:ASP:O	1:A:48:GLN:N	2.29	0.65
1:B:46:ASP:O	1:B:48:GLN:N	2.29	0.65
1:A:51:PHE:HE2	1:A:65:TYR:CB	2.09	0.65
1:A:188:LYS:O	1:A:190:VAL:N	2.30	0.65
1:B:60:CYS:O	1:B:64:CYS:SG	2.55	0.65
1:B:187:GLU:HG3	1:A:70:PRO:O	1.97	0.64
1:B:187:GLU:O	1:B:189:THR:N	2.30	0.63
1:A:39:ALA:HB2	1:A:191:PHE:CD1	2.32	0.63
1:B:65:TYR:O	1:B:67:HIS:N	2.32	0.62
1:A:190:VAL:O	1:A:192:THR:N	2.34	0.60
1:B:48:GLN:O	1:B:49:ALA:C	2.38	0.59

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:190:VAL:O	1:A:193:VAL:N	2.36	0.59
1:B:181:PHE:HE1	1:A:58:PRO:C	2.02	0.58
1:A:163:MET:HE2	1:A:163:MET:HA	1.85	0.58
1:A:163:MET:O	1:A:164:GLN:C	2.41	0.57
1:A:48:GLN:OE1	1:A:65:TYR:CD2	2.57	0.57
1:A:188:LYS:C	1:A:190:VAL:N	2.54	0.57
1:A:46:ASP:O	1:A:47:GLU:C	2.43	0.56
1:B:187:GLU:O	1:B:190:VAL:N	2.38	0.56
1:B:187:GLU:O	1:B:188:LYS:C	2.45	0.55
1:B:46:ASP:O	1:B:47:GLU:C	2.44	0.55
1:B:48:GLN:O	1:B:51:PHE:N	2.40	0.54
1:A:192:THR:O	1:A:194:PHE:N	2.42	0.53
1:A:51:PHE:CE2	1:A:65:TYR:CB	2.86	0.53
1:A:44:TRP:CE2	1:A:73:HIS:HB2	2.42	0.53
1:A:47:GLU:O	1:A:50:ASP:CA	2.57	0.52
1:A:191:PHE:O	1:A:192:THR:C	2.48	0.51
1:A:44:TRP:HZ2	1:A:73:HIS:CD2	2.27	0.51
1:B:65:TYR:CE2	1:B:69:PHE:HB3	2.46	0.51
1:A:189:THR:HG22	1:A:190:VAL:N	2.26	0.51
1:B:66:ASP:OD2	1:B:66:ASP:O	2.30	0.50
1:B:185:PRO:O	1:B:189:THR:OG1	2.30	0.50
1:A:47:GLU:O	1:A:48:GLN:O	2.30	0.49
1:B:48:GLN:O	1:B:50:ASP:N	2.44	0.49
1:A:48:GLN:C	1:A:50:ASP:N	2.66	0.49
1:B:181:PHE:CD1	1:A:59:GLY:CA	2.91	0.49
1:A:43:VAL:O	1:A:44:TRP:O	2.30	0.49
1:B:181:PHE:CE1	1:A:58:PRO:C	2.80	0.48
1:B:50:ASP:N	1:B:50:ASP:OD1	2.47	0.48
1:B:189:THR:O	1:B:192:THR:OG1	2.26	0.47
1:B:52:VAL:HG21	1:A:62:ASN:ND2	2.29	0.47
1:A:51:PHE:HE2	1:A:65:TYR:CA	2.28	0.47
1:A:189:THR:O	1:A:192:THR:OG1	2.30	0.47
1:A:18:THR:O	1:A:19:SER:OG	2.27	0.47
1:A:32:ARG:NH2	1:A:199:SER:OG	2.47	0.47
1:A:50:ASP:OD1	1:A:50:ASP:N	2.47	0.46
1:A:48:GLN:O	1:A:51:PHE:CB	2.63	0.46
1:A:184:ARG:N	1:A:185:PRO:HD3	2.30	0.46
1:B:47:GLU:O	1:B:48:GLN:C	2.54	0.45
1:A:193:VAL:HG12	1:A:194:PHE:CD1	2.51	0.45
1:B:188:LYS:O	1:B:192:THR:HG23	2.17	0.45
1:B:182:VAL:CG1	1:B:183:SER:N	2.80	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:71:ILE:O	1:A:158:TYR:OH	2.29	0.44
1:B:186:THR:O	1:B:187:GLU:C	2.54	0.44
1:B:183:SER:O	1:B:184:ARG:HB3	2.17	0.44
1:B:48:GLN:O	1:B:51:PHE:CB	2.66	0.44
1:B:65:TYR:HE2	1:B:69:PHE:HB3	1.82	0.44
1:A:191:PHE:O	1:A:192:THR:O	2.35	0.44
1:B:65:TYR:C	1:B:67:HIS:N	2.72	0.43
1:A:43:VAL:HG12	1:A:44:TRP:H	1.78	0.43
1:B:52:VAL:CG2	1:A:62:ASN:ND2	2.82	0.43
1:A:183:SER:O	1:A:184:ARG:O	2.37	0.42
1:A:77:TRP:CZ2	1:A:192:THR:CG2	2.99	0.42
1:A:181:PHE:N	1:A:181:PHE:CD2	2.87	0.42
1:B:162:SER:HA	1:B:189:THR:CG2	2.50	0.41
1:B:65:TYR:O	1:B:66:ASP:C	2.57	0.41
1:A:193:VAL:CG1	1:A:194:PHE:CD1	3.03	0.41
1:B:48:GLN:C	1:B:50:ASP:N	2.72	0.40
1:A:48:GLN:O	1:A:50:ASP:C	2.60	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:B:58:PRO:O	1:A:181:PHE:CE1[2_555]	2.07	0.13

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	156/226 (69%)	127 (81%)	15 (10%)	14 (9%)	1	16
1	B	151/226 (67%)	130 (86%)	15 (10%)	6 (4%)	4	38
All	All	307/452 (68%)	257 (84%)	30 (10%)	20 (6%)	1	25

All (20) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	47	GLU
1	A	47	GLU
1	A	48	GLN
1	A	185	PRO
1	A	193	VAL
1	B	54	ASN
1	B	188	LYS
1	A	43	VAL
1	A	54	ASN
1	A	192	THR
1	B	187	GLU
1	A	49	ALA
1	A	190	VAL
1	A	164	GLN
1	A	189	THR
1	B	48	GLN
1	B	66	ASP
1	A	44	TRP
1	A	184	ARG
1	A	95	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	144/204 (71%)	136 (94%)	8 (6%)	26 67
1	B	139/204 (68%)	133 (96%)	6 (4%)	35 74
All	All	283/408 (69%)	269 (95%)	14 (5%)	31 70

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	42	GLU
1	B	50	ASP
1	B	52	VAL

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	65	TYR
1	B	66	ASP
1	B	189	THR
1	A	44	TRP
1	A	50	ASP
1	A	163	MET
1	A	164	GLN
1	A	182	VAL
1	A	189	THR
1	A	192	THR
1	A	193	VAL

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

Mol	Chain	Res	Type
1	B	94	HIS
1	A	48	GLN

5.3.3 RNA

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates

There are no carbohydrates in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

5.7 Other polymers

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, 95^{th} 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	160/226 (70%)	0.00	3 (1%) 70 54	81, 125, 154, 178	0
1	B	155/226 (68%)	0.12	9 (5%) 26 16	89, 126, 160, 176	0
All	All	315/452 (69%)	0.06	12 (3%) 44 30	81, 126, 158, 178	0

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	24	TRP	4.8
1	A	93	MET	3.4
1	A	92	ALA	3.2
1	B	213	LEU	3.1
1	B	25	LEU	2.7
1	B	152	TYR	2.5
1	B	27	VAL	2.3
1	B	52	VAL	2.2
1	B	28	LEU	2.2
1	B	153	VAL	2.1
1	A	137	THR	2.0
1	B	91	VAL	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 carbohydrates in this entry.

6.4 Ligands

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.