



Full wwPDB X-ray Structure Validation Report ⓘ

May 23, 2020 – 12:02 am BST

PDB ID : 5LOC
Title : Crystal structure of the engineered D-Amino Acid Dehydrogenase (DAADH)
Authors : Dunstan, M.S.; Gahloth, D.
Deposited on : 2016-08-09
Resolution : 2.04 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 : 2.11
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) : 2.11

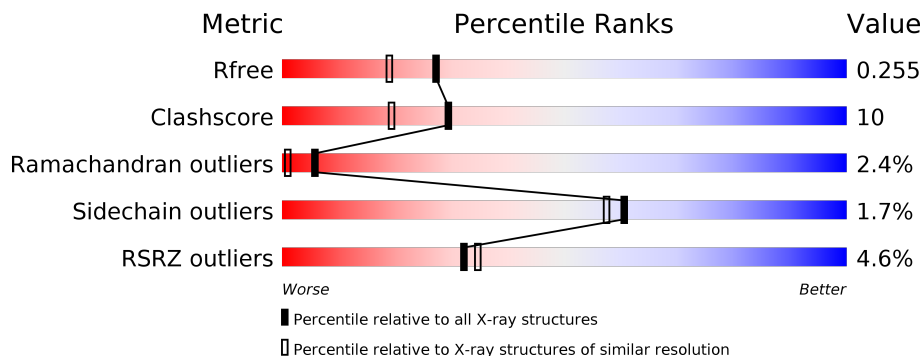
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.04 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	1692 (2.04-2.04)
Clashscore	141614	1773 (2.04-2.04)
Ramachandran outliers	138981	1752 (2.04-2.04)
Sidechain outliers	138945	1752 (2.04-2.04)
RSRZ outliers	127900	1672 (2.04-2.04)

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 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	320	
1	B	320	

2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 5126 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 Meso-diaminopimelate D-dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	320	2464	1540	437	476	11	0	0	0
1	B	317	2445	1530	433	471	11	0	0	0

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	2	GLY	-	expression tag	UNP P04964
A	151	LEU	GLN	conflict	UNP P04964
A	155	GLY	ASP	conflict	UNP P04964
A	170	ILE	THR	conflict	UNP P04964
A	175	ASP	ASP	conflict	UNP P04964
A	196	MET	ARG	conflict	UNP P04964
A	238	SER	SER	conflict	UNP P04964
A	245	ASN	HIS	conflict	UNP P04964
B	2	GLY	-	expression tag	UNP P04964
B	151	LEU	GLN	conflict	UNP P04964
B	155	GLY	ASP	conflict	UNP P04964
B	170	ILE	THR	conflict	UNP P04964
B	175	ASP	ASP	conflict	UNP P04964
B	196	MET	ARG	conflict	UNP P04964
B	238	SER	SER	conflict	UNP P04964
B	245	ASN	HIS	conflict	UNP P04964

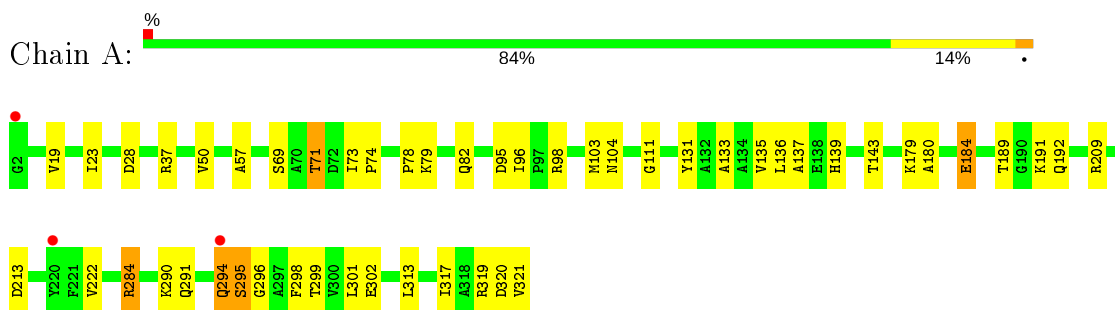
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	114	Total	O	0	0
			114	114		
2	B	103	Total	O	0	0
			103	103		

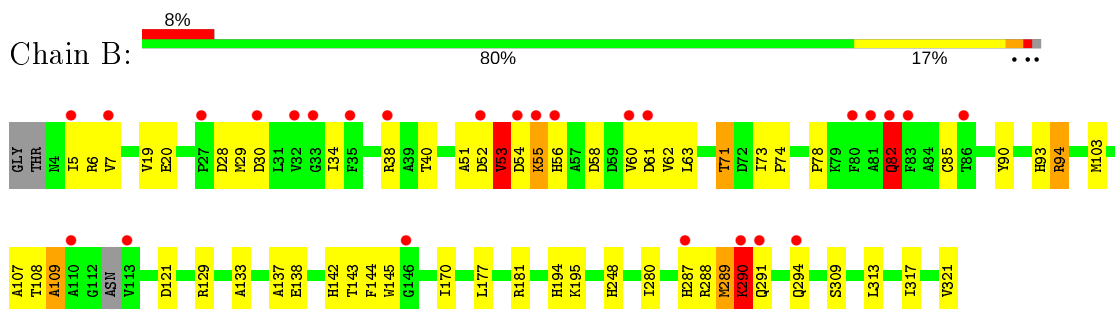
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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: Meso-diaminopimelate D-dehydrogenase



- Molecule 1: Meso-diaminopimelate D-dehydrogenase



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	120.53 Å 120.53 Å 97.63 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	71.30 – 2.04 71.30 – 2.04	Depositor EDS
% Data completeness (in resolution range)	100.0 (71.30-2.04) 100.0 (71.30-2.04)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.89 (at 2.03 Å)	Xtrriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.207 , 0.255 0.209 , 0.255	Depositor DCC
R_{free} test set	2549 reflections (4.86%)	wwPDB-VP
Wilson B-factor (Å ²)	44.6	Xtrriage
Anisotropy	0.029	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 54.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	0.025 for -h,-k,l	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	5126	wwPDB-VP
Average B, all atoms (Å ²)	66.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.20% of the height of the origin peak. No significant pseudotranslation is detected.*

¹ Intensities estimated from amplitudes.

² Theoretical 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	0.42	0/2516	0.60	1/3416 (0.0%)
1	B	0.47	0/2497	0.68	1/3390 (0.0%)
All	All	0.45	0/5013	0.64	2/6806 (0.0%)

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	2

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	B	290	LYS	CB-CG-CD	6.06	127.34	111.60
1	A	284	ARG	NE-CZ-NH1	-6.04	117.28	120.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	145	TRP	Peptide
1	B	53	VAL	Peptide

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	2464	0	2377	35	0
1	B	2445	0	2361	59	0
2	A	114	0	0	7	1
2	B	103	0	0	8	1
All	All	5126	0	4738	92	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 (92) 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:57:ALA:O	2:A:401:HOH:O	1.86	0.93
1:A:222:VAL:O	2:A:402:HOH:O	1.90	0.88
1:B:133:ALA:O	2:B:401:HOH:O	1.90	0.88
1:A:28:ASP:OD2	1:A:284:ARG:NH1	2.17	0.77
1:A:209:ARG:NH1	1:A:213:ASP:OD1	2.17	0.77
1:B:5:ILE:HG22	1:B:29:MET:HA	1.68	0.76
1:B:137:ALA:O	2:B:402:HOH:O	2.03	0.76
1:B:288:ARG:HA	1:B:290:LYS:HG2	1.68	0.76
1:B:142:HIS:ND1	2:B:405:HOH:O	2.19	0.75
1:A:299:THR:HG23	1:A:301:LEU:H	1.51	0.74
1:B:287:HIS:O	1:B:290:LYS:N	2.22	0.72
1:A:180:ALA:O	2:A:403:HOH:O	2.08	0.70
1:B:144:PHE:HB2	1:B:248:HIS:HB2	1.76	0.67
1:B:288:ARG:C	1:B:290:LYS:H	1.97	0.67
1:B:287:HIS:CE1	1:B:290:LYS:HD2	2.31	0.66
1:B:289:MET:O	2:B:403:HOH:O	2.15	0.64
1:B:54:ASP:OD1	1:B:55:LYS:N	2.31	0.64
1:B:61:ASP:OD1	1:B:62:VAL:N	2.31	0.64
1:B:93:HIS:O	1:B:94:ARG:HB3	1.97	0.63
1:B:61:ASP:O	1:B:85:CYS:N	2.32	0.62
1:B:6:ARG:HB2	1:B:61:ASP:OD1	2.01	0.61
1:B:317:ILE:O	2:B:404:HOH:O	2.16	0.61
1:B:288:ARG:HA	1:B:290:LYS:CG	2.31	0.61
1:B:5:ILE:HG21	1:B:29:MET:HG3	1.82	0.61
1:B:38:ARG:NH2	1:B:40:THR:O	2.36	0.59
1:B:288:ARG:NH1	1:B:309:SER:O	2.36	0.58
1:B:5:ILE:CG2	1:B:29:MET:HG3	2.34	0.58
1:B:19:VAL:HG13	1:B:280:ILE:HD11	1.84	0.58

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:139:HIS:HD2	2:A:404:HOH:O	1.87	0.57
1:B:288:ARG:CA	1:B:290:LYS:HG2	2.34	0.57
1:B:287:HIS:ND1	1:B:290:LYS:HD2	2.20	0.56
1:A:189:THR:HG22	1:A:192:GLN:HG3	1.90	0.54
1:A:104:ASN:OD1	1:A:296:GLY:HA3	2.08	0.54
1:A:111:GLY:O	1:A:290:LYS:NZ	2.41	0.53
1:B:287:HIS:ND1	1:B:290:LYS:CD	2.71	0.53
1:A:136:LEU:O	2:A:404:HOH:O	2.19	0.53
1:B:5:ILE:CG2	1:B:29:MET:HA	2.38	0.53
1:B:5:ILE:HD11	1:B:287:HIS:CG	2.44	0.52
1:B:51:ALA:O	1:B:53:VAL:HG12	2.09	0.52
1:A:299:THR:HG22	1:A:302:GLU:OE2	2.09	0.52
1:A:189:THR:HG23	1:A:191:LYS:H	1.74	0.52
1:A:19:VAL:O	1:A:23:ILE:HG12	2.11	0.51
1:B:61:ASP:CG	1:B:62:VAL:HG12	2.31	0.51
1:B:121:ASP:OD2	1:B:129:ARG:NH2	2.43	0.51
1:A:295:SER:HB2	1:A:298:PHE:HZ	1.77	0.50
1:B:288:ARG:C	1:B:290:LYS:N	2.64	0.49
1:A:319:ARG:O	1:A:321:VAL:N	2.47	0.48
1:A:37:ARG:HD3	1:A:50:VAL:HG21	1.95	0.48
1:A:73:ILE:HB	1:A:74:PRO:HD3	1.95	0.48
1:A:69:SER:OG	1:A:69:SER:O	2.29	0.48
1:B:291:GLN:OE1	1:B:291:GLN:N	2.47	0.48
1:A:133:ALA:O	2:A:404:HOH:O	2.19	0.48
1:B:7:VAL:HG22	1:B:30:ASP:O	2.14	0.48
1:A:95:ASP:OD1	1:A:98:ARG:HD3	2.14	0.48
1:A:28:ASP:CG	1:A:284:ARG:HH12	2.16	0.47
1:B:62:VAL:HB	1:B:85:CYS:HB2	1.97	0.47
1:A:299:THR:HG23	1:A:301:LEU:N	2.25	0.46
1:A:302:GLU:HG2	1:B:313:LEU:HD21	1.97	0.46
1:B:287:HIS:C	1:B:290:LYS:HG2	2.35	0.46
1:A:179:LYS:HE3	1:A:184:GLU:O	2.16	0.45
1:B:287:HIS:O	1:B:290:LYS:HG2	2.16	0.45
1:B:82:GLN:H	1:B:82:GLN:HG2	1.41	0.45
1:B:288:ARG:NH2	2:B:411:HOH:O	2.32	0.45
1:B:56:HIS:O	1:B:60:VAL:HG13	2.17	0.45
1:B:129:ARG:HH21	1:B:143:THR:HG21	1.82	0.44
1:A:96:ILE:HG21	1:B:321:VAL:HG21	2.00	0.44
1:B:61:ASP:OD2	1:B:62:VAL:HG12	2.16	0.44
1:B:321:VAL:HG22	2:B:404:HOH:O	2.18	0.44
1:A:73:ILE:HG12	1:A:73:ILE:H	1.69	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:6:ARG:HB2	1:B:61:ASP:CG	2.38	0.43
1:B:58:ASP:OD1	1:B:58:ASP:N	2.50	0.43
1:A:295:SER:HB2	1:A:298:PHE:CZ	2.53	0.43
1:A:313:LEU:O	1:A:317:ILE:HG13	2.18	0.43
1:B:170:ILE:HD11	1:B:194:HIS:HB3	2.00	0.43
1:B:287:HIS:O	1:B:289:MET:N	2.52	0.43
1:B:71:THR:C	1:B:74:PRO:HD2	2.39	0.43
1:B:78:PRO:N	1:B:103:MET:HE1	2.34	0.42
1:A:79:LYS:O	1:A:82:GLN:HG2	2.19	0.42
1:B:177:LEU:O	1:B:181:ARG:HG3	2.19	0.42
1:B:290:LYS:HE3	1:B:291:GLN:OE1	2.19	0.42
1:B:73:ILE:HG12	1:B:74:PRO:HD3	2.00	0.42
1:B:107:ALA:O	1:B:109:ALA:N	2.53	0.42
1:A:131:TYR:O	1:A:135:VAL:HG23	2.20	0.42
1:A:139:HIS:CD2	2:A:404:HOH:O	2.67	0.41
1:A:78:PRO:CD	1:A:103:MET:HE1	2.51	0.41
1:B:138:GLU:O	2:B:407:HOH:O	2.21	0.41
1:B:20:GLU:OE1	1:B:34:ILE:HD11	2.20	0.41
1:B:5:ILE:HG22	1:B:28:ASP:O	2.20	0.41
1:A:111:GLY:C	1:A:294:GLN:HG3	2.42	0.40
1:B:73:ILE:HD11	1:B:90:TYR:HB2	2.02	0.40
1:A:290:LYS:HA	1:A:294:GLN:HA	2.02	0.40
1:B:5:ILE:HD12	1:B:5:ILE:HA	1.86	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)
2:A:500:HOH:O	2:B:477:HOH:O[5_557]	2.12	0.08

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	318/320 (99%)	298 (94%)	14 (4%)	6 (2%)	8	2
1	B	315/320 (98%)	288 (91%)	18 (6%)	9 (3%)	4	0
All	All	633/640 (99%)	586 (93%)	32 (5%)	15 (2%)	6	1

All (15) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	137	ALA
1	A	143	THR
1	A	294	GLN
1	A	295	SER
1	B	53	VAL
1	B	109	ALA
1	A	71	THR
1	B	71	THR
1	B	108	THR
1	A	320	ASP
1	B	290	LYS
1	B	94	ARG
1	B	289	MET
1	B	82	GLN
1	B	294	GLN

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	259/259 (100%)	256 (99%)	3 (1%)	71	70
1	B	257/259 (99%)	251 (98%)	6 (2%)	50	44
All	All	516/518 (100%)	507 (98%)	9 (2%)	60	57

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	71	THR

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	184	GLU
1	A	291	GLN
1	B	52	ASP
1	B	55	LYS
1	B	63	LEU
1	B	82	GLN
1	B	195	LYS
1	B	290	LYS

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

Mol	Chain	Res	Type
1	A	294	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 carbohydrates 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	320/320 (100%)	-0.08	3 (0%) 84 86	39, 58, 94, 114	0
1	B	317/320 (99%)	0.31	26 (8%) 11 12	36, 65, 114, 143	0
All	All	637/640 (99%)	0.11	29 (4%) 32 35	36, 61, 108, 143	0

All (29) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	291	GLN	10.9
1	B	33	GLY	6.4
1	A	2	GLY	5.2
1	B	55	LYS	4.3
1	B	290	LYS	4.2
1	B	5	ILE	3.7
1	B	61	ASP	3.5
1	B	54	ASP	3.2
1	B	110	ALA	3.1
1	B	27	PRO	3.1
1	A	294	GLN	2.9
1	B	7	VAL	2.9
1	B	56	HIS	2.8
1	B	52	ASP	2.8
1	B	82	GLN	2.8
1	B	60	VAL	2.7
1	B	146	GLY	2.6
1	B	287	HIS	2.4
1	B	30	ASP	2.4
1	B	86	THR	2.4
1	A	220	TYR	2.4
1	B	35	PHE	2.3
1	B	83	PHE	2.3
1	B	113	VAL	2.3

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	B	294	GLN	2.2
1	B	81	ALA	2.1
1	B	32	VAL	2.1
1	B	80	PHE	2.1
1	B	38	ARG	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 carbohydrates 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.