



Full wwPDB X-ray Structure Validation Report ⓘ

Jun 12, 2024 – 06:21 AM EDT

PDB ID : 6NQ6
Title : Structure & function of a new Aspartylglucosaminuria variant
Authors : Pande, S.; Guo, H.C.
Deposited on : 2019-01-19
Resolution : 1.50 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	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.36.2
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)	:	2.36.2

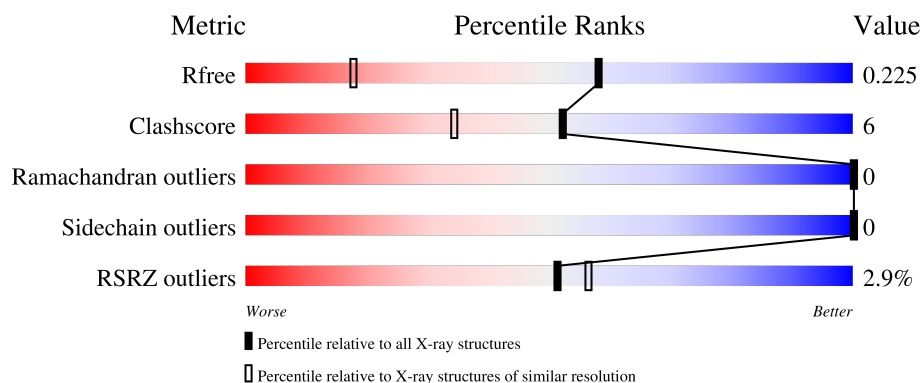
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.50 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	2936 (1.50-1.50)
Clashscore	141614	3144 (1.50-1.50)
Ramachandran outliers	138981	3066 (1.50-1.50)
Sidechain outliers	138945	3064 (1.50-1.50)
RSRZ outliers	127900	2884 (1.50-1.50)

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 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	138	<div> <div>4%</div> <div> <div></div> <div>85%</div> <div>12%</div> <div>..</div> </div> </div>
1	C	138	<div> <div>3%</div> <div> <div></div> <div>90%</div> <div>7%</div> <div>.</div> </div> </div>
2	B	144	<div> <div>2%</div> <div> <div></div> <div>88%</div> <div>10%</div> <div>..</div> </div> </div>
2	D	144	<div> <div>3%</div> <div> <div></div> <div>85%</div> <div>12%</div> <div>..</div> </div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4421 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 N(4)-(Beta-N-acetylglucosaminy)-L-asparaginase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	134	Total	C	N	O	S	0	0	0
			1039	654	180	199	6			
1	C	134	Total	C	N	O	S	0	0	0
			1039	654	180	199	6			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	99	LYS	THR	engineered mutation	UNP A0A376EJJ1
C	99	LYS	THR	engineered mutation	UNP A0A376EJJ1

- Molecule 2 is a protein called N(4)-(Beta-N-acetylglucosaminy)-L-asparaginase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	144	Total	C	N	O	S	0	2	0
			1092	679	201	205	7			
2	D	144	Total	C	N	O	S	0	2	0
			1089	678	200	204	7			

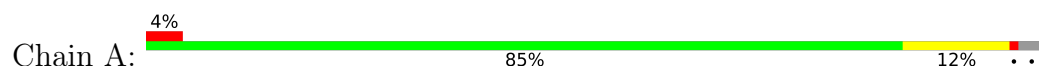
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	41	Total	O	0	0
			41	41		
3	B	41	Total	O	0	0
			41	41		
3	C	49	Total	O	0	0
			49	49		
3	D	31	Total	O	0	0
			31	31		

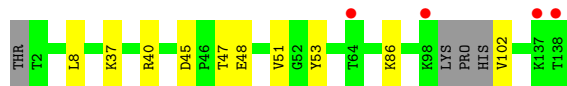
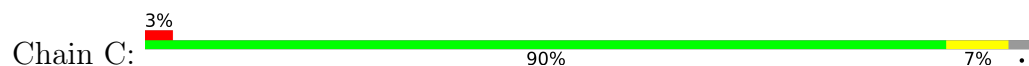
3 Residue-property plots [i](#)

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.

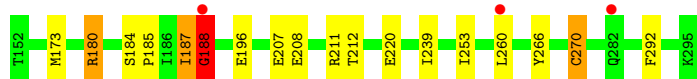
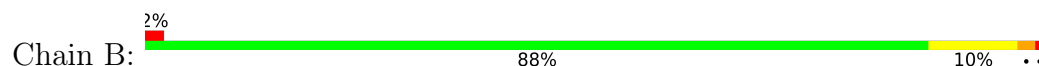
- Molecule 1: N(4)-(Beta-N-acetylglucosaminy)-L-asparaginase



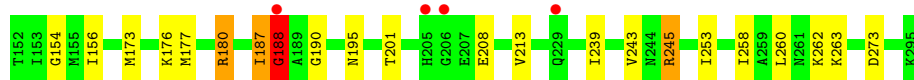
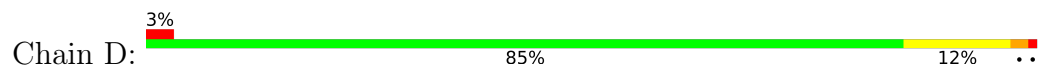
- Molecule 1: N(4)-(Beta-N-acetylglucosaminy)-L-asparaginase



- Molecule 2: N(4)-(Beta-N-acetylglucosaminy)-L-asparaginase



- Molecule 2: N(4)-(Beta-N-acetylglucosaminy)-L-asparaginase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	46.15Å 96.02Å 61.49Å 90.00° 90.05° 90.00°	Depositor
Resolution (Å)	19.99 – 1.50 19.99 – 1.50	Depositor EDS
% Data completeness (in resolution range)	99.0 (19.99-1.50) 99.0 (19.99-1.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.91 (at 1.50Å)	Xtriage
Refinement program	REFMAC 5.8.0189	Depositor
R, R_{free}	0.194 , 0.224 0.198 , 0.225	Depositor DCC
R_{free} test set	4333 reflections (5.12%)	wwPDB-VP
Wilson B-factor (Å ²)	11.3	Xtriage
Anisotropy	0.046	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.47 , 49.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.043 for h,-k,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4421	wwPDB-VP
Average B, all atoms (Å ²)	14.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.71% 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	1.10	5/1056 (0.5%)	1.14	8/1423 (0.6%)
1	C	0.99	1/1056 (0.1%)	1.00	2/1423 (0.1%)
2	B	1.21	5/1109 (0.5%)	1.19	6/1490 (0.4%)
2	D	1.21	5/1109 (0.5%)	1.24	9/1490 (0.6%)
All	All	1.13	16/4330 (0.4%)	1.15	25/5826 (0.4%)

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
2	B	0	2
2	D	0	1
All	All	0	4

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	188	GLY	N-CA	16.59	1.71	1.46
2	B	188	GLY	N-CA	14.09	1.67	1.46
2	B	188	GLY	CA-C	11.60	1.70	1.51
2	D	188	GLY	C-O	11.42	1.42	1.23
1	A	54	GLY	N-CA	11.05	1.62	1.46
2	B	270	CYS	CB-SG	-8.56	1.67	1.82
2	D	188	GLY	CA-C	7.74	1.64	1.51
1	A	54	GLY	CA-C	7.72	1.64	1.51
2	B	188	GLY	C-O	7.70	1.35	1.23
1	A	102	VAL	N-CA	-7.20	1.31	1.46
1	A	20	GLU	CD-OE1	-6.46	1.18	1.25
1	C	48	GLU	CD-OE2	6.25	1.32	1.25
1	A	83	GLU	CD-OE2	-5.51	1.19	1.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	220	GLU	CD-OE2	-5.35	1.19	1.25
2	D	180[A]	ARG	CA-CB	-5.27	1.42	1.53
2	D	180[B]	ARG	CA-CB	-5.27	1.42	1.53

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	187	ILE	C-N-CA	-12.86	95.29	122.30
2	B	187	ILE	C-N-CA	-11.30	98.58	122.30
2	D	180[A]	ARG	NE-CZ-NH1	-10.36	115.12	120.30
2	D	180[B]	ARG	NE-CZ-NH1	-10.36	115.12	120.30
2	B	180[A]	ARG	NE-CZ-NH1	-9.31	115.65	120.30
2	B	180[B]	ARG	NE-CZ-NH1	-9.31	115.65	120.30
1	C	40	ARG	NE-CZ-NH1	8.48	124.54	120.30
1	A	40	ARG	NE-CZ-NH2	-7.78	116.41	120.30
1	A	40	ARG	NE-CZ-NH1	7.58	124.09	120.30
1	A	53	TYR	C-N-CA	-7.50	106.55	122.30
2	B	188	GLY	N-CA-C	7.22	131.15	113.10
1	A	93	ARG	NE-CZ-NH2	-7.16	116.72	120.30
2	D	245	ARG	NE-CZ-NH1	6.69	123.65	120.30
2	D	187	ILE	O-C-N	-6.62	111.95	123.20
1	A	56	ARG	NE-CZ-NH1	6.51	123.56	120.30
1	A	8	LEU	CA-CB-CG	6.21	129.57	115.30
1	A	86	LYS	CD-CE-NZ	6.11	125.75	111.70
2	D	260	LEU	CA-CB-CG	5.99	129.07	115.30
1	C	40	ARG	NE-CZ-NH2	-5.78	117.41	120.30
2	D	188	GLY	N-CA-C	5.59	127.08	113.10
1	A	54	GLY	N-CA-C	5.46	126.76	113.10
2	B	211	ARG	NE-CZ-NH1	5.37	122.98	120.30
2	D	187	ILE	CA-C-N	5.12	126.44	116.20
2	B	187	ILE	O-C-N	-5.07	114.58	123.20
2	D	213	VAL	CG1-CB-CG2	-5.03	102.85	110.90

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	54	GLY	Peptide
2	B	188	GLY	Mainchain,Peptide
2	D	188	GLY	Peptide

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	1039	0	1037	8	0
1	C	1039	0	1037	6	0
2	B	1092	0	1100	16	0
2	D	1089	0	1101	25	0
3	A	41	0	0	1	0
3	B	41	0	0	0	0
3	C	49	0	0	2	0
3	D	31	0	0	1	0
All	All	4421	0	4275	47	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 (47) 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)
2:D:188:GLY:N	2:D:188:GLY:CA	1.70	1.48
1:A:74:TYR:OH	2:D:245:ARG:NH2	1.74	1.19
2:D:187:ILE:O	2:D:188:GLY:CA	1.97	1.08
2:D:187:ILE:C	2:D:188:GLY:CA	2.22	1.07
2:D:187:ILE:O	2:D:188:GLY:HA2	1.82	0.80
2:B:260:LEU:HD12	2:B:266:TYR:HB3	1.66	0.77
1:C:45:ASP:OD1	1:C:47:THR:OG1	2.04	0.76
1:A:6:ILE:HG23	2:B:260:LEU:HD21	1.68	0.74
2:D:273:ASP:OD1	3:D:301:HOH:O	2.06	0.73
2:D:195:ASN:O	2:D:262:LYS:HD3	1.94	0.68
2:B:207[A]:GLU:HG3	1:C:102:VAL:HG22	1.79	0.64
2:D:239:ILE:O	2:D:243:VAL:HG23	1.97	0.63
1:A:74:TYR:CZ	2:D:245:ARG:NH2	2.67	0.61
2:D:154:GLY:HA3	2:D:201:THR:HG21	1.88	0.54
2:D:208:GLU:HG3	2:D:253:ILE:HG21	1.89	0.54
1:A:48:GLU:O	1:A:54:GLY:HA3	2.09	0.53
2:D:208:GLU:HG3	2:D:253:ILE:CG2	2.39	0.52
2:B:196:GLU:HG2	2:D:245:ARG:NH2	2.25	0.52
2:D:187:ILE:O	2:D:188:GLY:C	2.48	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:6:ILE:CG2	2:B:260:LEU:HD21	2.39	0.50
2:D:156:ILE:HG23	2:D:201:THR:OG1	2.12	0.49
1:A:80:ALA:HB3	1:A:105:VAL:HG12	1.94	0.49
2:B:212:THR:HG21	2:B:239:ILE:HG12	1.95	0.48
2:B:208:GLU:HG3	2:B:253:ILE:HG23	1.96	0.48
2:B:207[A]:GLU:H	2:B:207[A]:GLU:CD	2.17	0.48
3:A:218:HOH:O	2:D:177:MET:HG2	2.13	0.48
1:C:53:TYR:HB2	1:C:86:LYS:HG3	1.97	0.47
2:D:173:MET:H	2:D:180[A]:ARG:NH1	2.14	0.46
2:B:207[A]:GLU:CD	2:B:207[A]:GLU:N	2.69	0.46
1:C:8:LEU:HB3	2:D:258:ILE:HD12	1.98	0.46
2:B:260:LEU:CD1	2:B:266:TYR:HB3	2.42	0.46
1:C:51:VAL:O	3:C:201:HOH:O	2.21	0.46
1:A:131:GLU:HA	1:A:134:GLU:HG3	1.98	0.44
2:D:190:GLY:O	2:D:201:THR:HG23	2.17	0.44
2:D:263:LYS:HA	2:D:263:LYS:HD3	1.80	0.44
2:B:173:MET:H	2:B:180[A]:ARG:NH1	2.16	0.44
1:A:6:ILE:HD13	2:B:260:LEU:HD11	1.98	0.43
2:D:176:LYS:HB2	2:D:180[B]:ARG:HD3	2.00	0.43
2:D:195:ASN:O	2:D:262:LYS:CD	2.64	0.43
1:C:37:LYS:HD3	3:C:239:HOH:O	2.18	0.43
2:B:187:ILE:O	2:B:188:GLY:C	2.50	0.42
2:B:184:SER:N	2:B:185:PRO:CD	2.82	0.42
2:B:270:CYS:O	2:B:292:PHE:HA	2.20	0.42
2:D:177:MET:HE3	2:D:177:MET:HB3	1.88	0.42
2:D:195:ASN:O	2:D:262:LYS:CE	2.67	0.42

There are no symmetry-related clashes.

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	130/138 (94%)	126 (97%)	4 (3%)	0	100	100
1	C	130/138 (94%)	128 (98%)	2 (2%)	0	100	100
2	B	144/144 (100%)	139 (96%)	5 (4%)	0	100	100
2	D	144/144 (100%)	138 (96%)	6 (4%)	0	100	100
All	All	548/564 (97%)	531 (97%)	17 (3%)	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	111/115 (96%)	111 (100%)	0	100	100
1	C	111/115 (96%)	111 (100%)	0	100	100
2	B	112/110 (102%)	112 (100%)	0	100	100
2	D	112/110 (102%)	112 (100%)	0	100	100
All	All	446/450 (99%)	446 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 [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	134/138 (97%)	0.11	5 (3%) 41 46	6, 12, 28, 47	0
1	C	134/138 (97%)	0.15	4 (2%) 50 55	7, 13, 32, 48	0
2	B	144/144 (100%)	0.02	3 (2%) 63 68	6, 12, 19, 33	0
2	D	144/144 (100%)	0.06	4 (2%) 53 57	6, 12, 24, 31	0
All	All	556/564 (98%)	0.08	16 (2%) 51 56	6, 12, 26, 48	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	64	THR	6.7
2	B	260	LEU	4.9
2	D	188	GLY	3.7
1	C	138	THR	3.6
1	C	137	LYS	3.3
1	A	54	GLY	2.9
2	D	229	GLN	2.7
2	D	205	HIS	2.7
2	B	188	GLY	2.7
1	C	98	LYS	2.6
1	A	137	LYS	2.6
1	A	134	GLU	2.5
2	D	206	GLY	2.4
2	B	282	GLN	2.3
1	A	2	THR	2.1
1	A	74	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 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.