



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 18, 2026 – 10:43 AM UTC

PDB ID : 1CRW / pdb_00001crw
Title : CRYSTAL STRUCTURE OF APO-GLYCERALDEHYDE-3-PHOSPHAT
E DEHYDROGENASE FROM PALINURUS VERSICOLOR AT 2.0A RES-
OLUTION
Authors : Shen, Y.; Li, J.; Song, S.; Lin, Z.
Deposited on : 1999-08-16
Resolution : 2.00 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

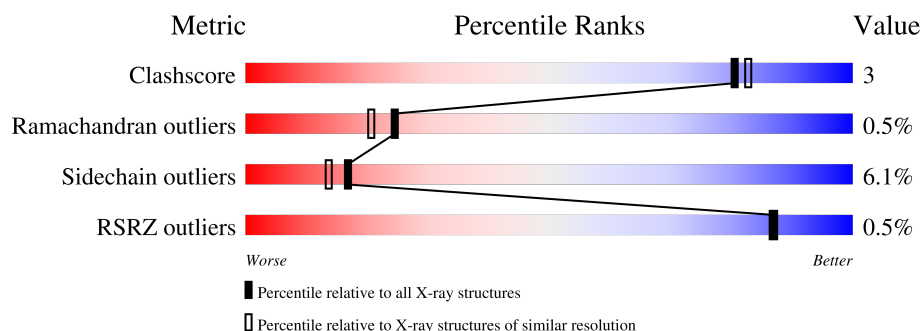
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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))
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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	G	333	
1	R	333	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 5337 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 D-GLYCERALDEHYDE-3-PHOSPHATE-DEHYDROGEN ASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	G	333	Total	C	N	O	S	0	0	0
			2506	1589	421	480	16			
1	R	333	Total	C	N	O	S	0	0	0
			2506	1589	421	480	16			

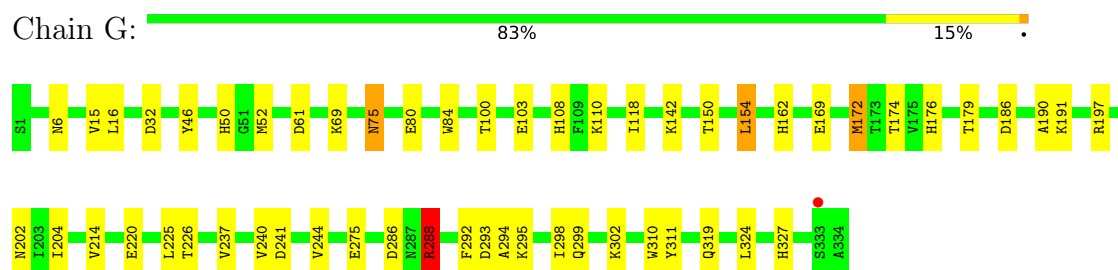
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	G	194	Total	O	0	0
			194	194		
2	R	131	Total	O	0	0
			131	131		

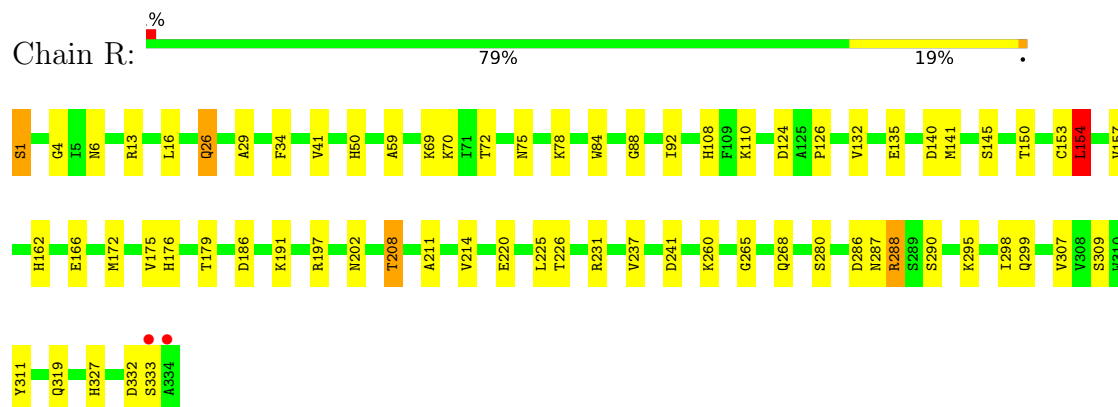
3 Residue-property plots

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: D-GLYCERALDEHYDE-3-PHOSPHATE-DEHYDROGENASE



• Molecule 1: D-GLYCERALDEHYDE-3-PHOSPHATE-DEHYDROGENASE



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	128.45Å 99.84Å 80.79Å 90.00° 115.14° 90.00°	Depositor
Resolution (Å)	6.00 – 2.00 6.00 – 2.00	Depositor EDS
% Data completeness (in resolution range)	74.5 (6.00-2.00) 66.9 (6.00-2.00)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.46 (at 1.80Å)	Xtriage
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.165 , 0.226 0.161 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	10.4	Xtriage
Anisotropy	0.819	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 87.1	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	0.95	EDS
Total number of atoms	5337	wwPDB-VP
Average B, all atoms (Å ²)	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.53% 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	G	1.03	5/2551 (0.2%)	1.67	28/3446 (0.8%)
1	R	0.98	6/2551 (0.2%)	1.70	37/3446 (1.1%)
All	All	1.01	11/5102 (0.2%)	1.69	65/6892 (0.9%)

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	G	50	HIS	CD2-NE2	-6.97	1.30	1.37
1	R	176	HIS	CD2-NE2	-6.86	1.30	1.37
1	R	327	HIS	CD2-NE2	-6.69	1.30	1.37
1	G	108	HIS	CD2-NE2	-6.65	1.30	1.37
1	R	108	HIS	CD2-NE2	-6.38	1.30	1.37
1	G	162	HIS	CD2-NE2	-6.31	1.30	1.37
1	G	176	HIS	CD2-NE2	-6.27	1.30	1.37
1	R	50	HIS	CD2-NE2	-6.15	1.31	1.37
1	R	41	VAL	CA-CB	5.54	1.60	1.54
1	R	162	HIS	CD2-NE2	-5.50	1.31	1.37
1	G	50	HIS	CG-ND1	-5.01	1.32	1.38

All (65) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	R	186	ASP	CA-CB-CG	8.49	121.09	112.60
1	R	6	ASN	OD1-CG-ND2	-8.16	114.44	122.60
1	G	226	THR	CA-CB-OG1	-7.50	98.34	109.60
1	R	226	THR	CA-CB-OG1	-7.36	98.56	109.60
1	G	6	ASN	OD1-CG-ND2	-7.27	115.33	122.60
1	R	150	THR	CA-C-O	-7.11	112.88	120.42
1	G	75	ASN	OD1-CG-ND2	-7.11	115.49	122.60
1	R	202	ASN	N-CA-C	6.94	120.47	109.50
1	G	197	ARG	NE-CZ-NH2	6.57	125.11	119.20
1	R	13	ARG	N-CA-CB	-6.56	100.48	110.12
1	R	176	HIS	CB-CG-CD2	-6.35	122.94	131.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	R	299	GLN	CA-CB-CG	-6.27	101.56	114.10
1	R	6	ASN	CB-CG-ND2	6.23	125.75	116.40
1	G	162	HIS	CB-CG-CD2	-6.18	123.17	131.20
1	R	124	ASP	CA-CB-CG	6.14	118.74	112.60
1	R	34	PHE	N-CA-C	6.07	120.85	112.90
1	R	268	GLN	OE1-CD-NE2	-6.04	116.56	122.60
1	G	84	TRP	CG-CD2-CE3	5.99	139.89	133.90
1	G	176	HIS	CB-CG-CD2	-5.98	123.42	131.20
1	R	4	GLY	O-C-N	-5.97	116.92	123.49
1	G	244	VAL	N-CA-CB	-5.90	101.77	111.45
1	G	186	ASP	CA-CB-CG	5.89	118.49	112.60
1	R	208	THR	O-C-N	-5.88	116.77	123.48
1	G	108	HIS	CB-CG-CD2	-5.87	123.57	131.20
1	G	310	TRP	CE2-CD2-CG	-5.83	100.20	107.20
1	G	80	GLU	CA-CB-CG	5.79	125.68	114.10
1	G	288	ARG	NE-CZ-NH1	5.78	127.28	121.50
1	G	15	VAL	N-CA-C	-5.76	104.90	110.72
1	R	78	LYS	CA-C-N	5.57	125.67	119.32
1	R	78	LYS	C-N-CA	5.57	125.67	119.32
1	R	26	GLN	OE1-CD-NE2	-5.56	117.04	122.60
1	G	204	ILE	CA-C-O	-5.55	117.11	120.88
1	G	6	ASN	CB-CG-ND2	5.51	124.67	116.40
1	G	202	ASN	N-CA-C	5.47	118.38	109.24
1	R	197	ARG	NE-CZ-NH2	5.46	124.12	119.20
1	G	324	LEU	N-CA-C	-5.46	105.25	111.14
1	R	295	LYS	CB-CG-CD	-5.38	98.92	111.30
1	R	75	ASN	OD1-CG-ND2	-5.36	117.24	122.60
1	G	226	THR	CA-CB-CG2	5.36	119.61	110.50
1	R	327	HIS	CB-CG-CD2	-5.36	124.23	131.20
1	R	280	SER	N-CA-C	5.35	117.86	111.71
1	G	299	GLN	OE1-CD-NE2	-5.31	117.29	122.60
1	R	265	GLY	O-C-N	-5.29	116.48	121.77
1	R	154	LEU	CB-CA-C	-5.29	102.55	110.90
1	R	132	VAL	N-CA-CB	-5.28	106.34	111.57
1	R	287	ASN	OD1-CG-ND2	-5.28	117.32	122.60
1	R	260	LYS	CB-CG-CD	-5.27	99.17	111.30
1	G	327	HIS	CB-CG-CD2	-5.20	124.44	131.20
1	R	132	VAL	O-C-N	-5.19	117.91	121.85
1	R	332	ASP	CA-CB-CG	5.18	117.78	112.60
1	R	226	THR	CA-CB-CG2	5.18	119.30	110.50
1	R	175	VAL	N-CA-C	-5.16	99.67	107.73
1	R	140	ASP	CA-CB-CG	5.15	117.75	112.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	84	TRP	CB-CG-CD1	-5.10	119.25	126.90
1	G	190	ALA	N-CA-C	5.09	117.49	111.33
1	R	287	ASN	CA-CB-CG	5.09	117.69	112.60
1	R	241	ASP	CA-CB-CG	5.08	117.68	112.60
1	G	293	ASP	CA-CB-CG	5.07	117.67	112.60
1	R	26	GLN	O-C-N	5.07	129.20	123.27
1	G	288	ARG	NE-CZ-NH2	-5.07	114.64	119.20
1	R	59	ALA	N-CA-C	-5.06	101.47	109.76
1	R	84	TRP	CE2-CD2-CG	-5.04	101.16	107.20
1	G	100	THR	N-CA-C	5.02	119.55	113.38
1	G	310	TRP	CG-CD2-CE3	5.02	138.92	133.90
1	G	241	ASP	CA-CB-CG	5.00	117.60	112.60

There are no chirality outliers.

There are no planarity outliers.

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	G	2506	0	2529	13	0
1	R	2506	0	2529	13	0
2	G	194	0	0	1	0
2	R	131	0	0	2	0
All	All	5337	0	5058	26	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 3.

All (26) 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:G:154:LEU:HD13	1:G:214:VAL:HG21	1.68	0.74
1:R:154:LEU:HD13	1:R:214:VAL:HG21	1.71	0.71
1:G:154:LEU:HG	1:G:172:MET:HE3	1.87	0.57
1:G:286:ASP:OD2	1:G:288:ARG:HD3	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:R:126:PRO:HG3	1:R:141:MET:HE2	1.93	0.50
1:R:1:SER:N	1:R:26:GLN:HB3	2.26	0.50
1:R:211:ALA:HB1	1:R:225:LEU:HB3	1.93	0.50
1:R:286:ASP:OD2	1:R:288:ARG:HD3	2.11	0.48
1:G:172:MET:HE2	1:G:240:VAL:HG22	1.93	0.48
1:R:220:GLU:HB2	2:R:421:HOH:O	2.14	0.48
1:G:150:THR:HG22	1:G:172:MET:HE1	1.97	0.46
1:G:46:TYR:HE2	1:G:52:MET:HE2	1.81	0.45
1:R:69:LYS:HD2	1:R:69:LYS:H	1.82	0.44
1:R:29:ALA:HA	1:R:72:THR:O	2.17	0.44
1:R:172:MET:HE1	1:R:208:THR:HG21	1.99	0.43
1:R:110:LYS:HD3	1:R:110:LYS:HA	1.75	0.43
1:G:191:LYS:HD3	1:G:191:LYS:HA	1.86	0.42
1:G:32:ASP:O	1:G:75:ASN:HA	2.20	0.41
1:G:150:THR:CG2	1:G:172:MET:HE1	2.50	0.41
1:G:292:PHE:CE2	1:G:294:ALA:HA	2.56	0.41
1:R:179:THR:OG1	1:R:231:ARG:NH2	2.54	0.41
1:G:179:THR:HB	2:G:460:HOH:O	2.20	0.41
1:G:275:GLU:O	1:G:295:LYS:HD3	2.21	0.41
1:R:88:GLY:HA3	2:R:387:HOH:O	2.21	0.40
1:R:153:CYS:HA	1:R:290:SER:HB2	2.02	0.40
1:G:142:LYS:HB3	1:G:142:LYS:HE2	1.88	0.40

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	G	331/333 (99%)	319 (96%)	11 (3%)	1 (0%)	36	35
1	R	331/333 (99%)	319 (96%)	10 (3%)	2 (1%)	21	17
All	All	662/666 (99%)	638 (96%)	21 (3%)	3 (0%)	24	21

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	G	237	VAL
1	R	237	VAL
1	R	166	GLU

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	G	269/269 (100%)	252 (94%)	17 (6%)	16	13
1	R	269/269 (100%)	253 (94%)	16 (6%)	18	14
All	All	538/538 (100%)	505 (94%)	33 (6%)	17	14

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	G	16	LEU
1	G	61	ASP
1	G	69	LYS
1	G	103	GLU
1	G	110	LYS
1	G	118	ILE
1	G	154	LEU
1	G	169	GLU
1	G	172	MET
1	G	174	THR
1	G	220	GLU
1	G	225	LEU
1	G	288	ARG
1	G	298	ILE
1	G	302	LYS
1	G	311	TYR
1	G	319	GLN
1	R	1	SER
1	R	16	LEU
1	R	70	LYS

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Mol	Chain	Res	Type
1	R	92	ILE
1	R	135	GLU
1	R	145	SER
1	R	154	LEU
1	R	157	VAL
1	R	191	LYS
1	R	288	ARG
1	R	298	ILE
1	R	307	VAL
1	R	309	SER
1	R	311	TYR
1	R	319	GLN
1	R	333	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	G	268	GLN
1	G	287	ASN
1	R	75	ASN
1	R	81	ASN
1	R	146	ASN

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 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 [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	G	333/333 (100%)	-1.15	1 (0%) 90 89	6, 17, 35, 66	0
1	R	333/333 (100%)	-0.90	2 (0%) 85 85	8, 23, 51, 82	0
All	All	666/666 (100%)	-1.03	3 (0%) 87 87	6, 19, 48, 82	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	R	333	SER	2.7
1	G	333	SER	2.6
1	R	334	ALA	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.