



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 04:09 AM UTC

PDB ID : 5O0R / pdb_00005o0r
Title : Deglycosylated Nogo Receptor with native disulfide structure
Authors : Pronker, M.F.; Janssen, B.J.C.
Deposited on : 2017-05-16
Resolution : 2.50 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 : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
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

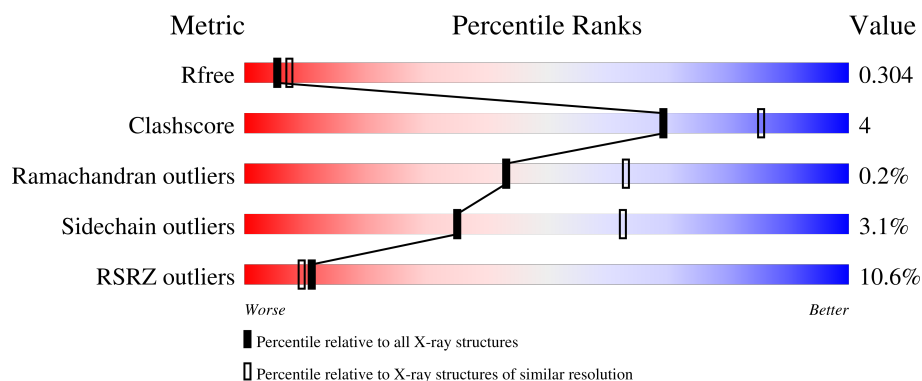
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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}	180053	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	334	 9% 78% 7% 13%
1	B	334	 10% 75% 10% 15%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 4619 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 Reticulon-4 receptor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	289	Total	C	N	O	S	0	0	0
			2269	1433	425	396	15			
1	B	284	Total	C	N	O	S	0	0	0
			2236	1413	419	391	13			

There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	24	GLY	-	expression tag	UNP Q99PI8
A	25	SER	-	expression tag	UNP Q99PI8
A	349	ALA	-	expression tag	UNP Q99PI8
A	350	ALA	-	expression tag	UNP Q99PI8
A	351	ALA	-	expression tag	UNP Q99PI8
A	352	HIS	-	expression tag	UNP Q99PI8
A	353	HIS	-	expression tag	UNP Q99PI8
A	354	HIS	-	expression tag	UNP Q99PI8
A	355	HIS	-	expression tag	UNP Q99PI8
A	356	HIS	-	expression tag	UNP Q99PI8
A	357	HIS	-	expression tag	UNP Q99PI8
B	24	GLY	-	expression tag	UNP Q99PI8
B	25	SER	-	expression tag	UNP Q99PI8
B	349	ALA	-	expression tag	UNP Q99PI8
B	350	ALA	-	expression tag	UNP Q99PI8
B	351	ALA	-	expression tag	UNP Q99PI8
B	352	HIS	-	expression tag	UNP Q99PI8
B	353	HIS	-	expression tag	UNP Q99PI8
B	354	HIS	-	expression tag	UNP Q99PI8
B	355	HIS	-	expression tag	UNP Q99PI8
B	356	HIS	-	expression tag	UNP Q99PI8
B	357	HIS	-	expression tag	UNP Q99PI8

- Molecule 2 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	B	1	Total	C	N	O	0	0
			14	8	1	5		
2	B	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		

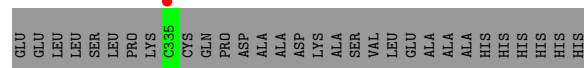
- Molecule 4 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Cl	0	0
			1	1		
4	B	2	Total	Cl	0	0
			2	2		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	20	Total	O	0	0
			20	20		
5	B	9	Total	O	0	0
			9	9		

- Molecule 1: Reticulon-4 receptor



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	46.73Å 111.65Å 114.83Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	43.10 – 2.50 43.10 – 2.50	Depositor EDS
% Data completeness (in resolution range)	98.4 (43.10-2.50) 98.4 (43.10-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.21	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.53 (at 2.51Å)	Xtriage
Refinement program	PHENIX (1.11.1_2575: ???)	Depositor
R, R_{free}	0.252 , 0.303 0.252 , 0.304	Depositor DCC
R_{free} test set	984 reflections (4.08%)	wwPDB-VP
Wilson B-factor (Å ²)	39.1	Xtriage
Anisotropy	0.609	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 50.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.000 for -h,l,k	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	4619	wwPDB-VP
Average B, all atoms (Å ²)	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 35.03 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 6.2171e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: CL, NAG, GOL

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	0.10	0/2327	0.31	0/3167
1	B	0.10	0/2293	0.30	0/3122
All	All	0.10	0/4620	0.31	0/6289

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	2269	0	2244	16	0
1	B	2236	0	2211	16	0
2	A	42	0	39	0	0
2	B	28	0	26	2	0
3	A	6	0	8	0	0
3	B	6	0	8	0	0
4	A	1	0	0	0	0
4	B	2	0	0	0	0
5	A	20	0	0	0	0
5	B	9	0	0	0	0
All	All	4619	0	4536	32	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 4.

All (32) 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:61:ARG:HG2	1:A:85:ILE:HB	1.83	0.60
1:A:219:PRO:HA	1:A:245:VAL:HG23	1.86	0.56
1:A:36:GLU:HG3	1:A:37:PRO:HA	1.88	0.55
1:A:167:GLN:OE1	1:A:189:ARG:NH2	2.42	0.53
1:B:38:LYS:HB2	1:B:61:ARG:HD3	1.90	0.53
1:A:87:TRP:HA	1:A:111:ASP:HB3	1.91	0.53
1:B:64:LEU:HB2	1:B:88:LEU:HD23	1.91	0.52
1:A:244:GLU:N	1:A:244:GLU:OE2	2.43	0.51
1:A:252:LEU:HD21	1:A:255:LEU:HB2	1.93	0.50
1:B:84:THR:HA	1:B:107:LEU:HA	1.92	0.49
1:B:290:PRO:HD2	1:B:293:LEU:HD12	1.95	0.49
1:B:111:ASP:HA	1:B:136:HIS:HB2	1.96	0.48
1:B:233:LEU:O	1:B:236:ASN:ND2	2.47	0.47
1:A:78:GLN:O	1:A:79:SER:OG	2.30	0.47
1:A:194:PRO:HG2	1:A:197:ALA:HB2	1.98	0.46
1:B:81:ARG:HB2	2:B:401:NAG:H82	1.98	0.46
1:B:272:TRP:O	1:B:276:GLN:HG2	2.17	0.45
1:A:150:PHE:HB3	1:A:177:LEU:HD21	1.98	0.45
1:A:157:GLN:HA	1:A:180:LEU:HA	1.97	0.45
1:B:194:PRO:HD2	1:B:197:ALA:HB2	1.99	0.44
1:A:138:ASP:OD1	1:A:138:ASP:N	2.43	0.44
1:A:260:ASN:HB2	1:A:262:TRP:NE1	2.33	0.44
1:B:157:GLN:HA	1:B:180:LEU:HA	2.00	0.44
1:A:61:ARG:HD2	1:A:85:ILE:HD12	1.99	0.43
1:B:177:LEU:HB3	1:B:180:LEU:HB2	2.00	0.43
1:B:258:ASN:O	1:B:260:ASN:ND2	2.52	0.43
1:B:291:GLN:H	1:B:291:GLN:CD	2.26	0.43
1:A:64:LEU:HB2	1:A:88:LEU:HD23	2.00	0.43
1:B:155:ALA:HB1	2:B:402:NAG:H82	2.01	0.42
1:A:272:TRP:O	1:A:276:GLN:HG2	2.21	0.41
1:B:285:VAL:HB	1:B:298:LEU:HB3	2.03	0.41
1:B:234:PHE:HA	1:B:260:ASN:HD21	1.86	0.41

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	285/334 (85%)	262 (92%)	23 (8%)	0	100	100
1	B	281/334 (84%)	259 (92%)	21 (8%)	1 (0%)	30	49
All	All	566/668 (85%)	521 (92%)	44 (8%)	1 (0%)	43	63

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	75	ALA

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	247/283 (87%)	239 (97%)	8 (3%)	34	62
1	B	243/283 (86%)	236 (97%)	7 (3%)	37	65
All	All	490/566 (87%)	475 (97%)	15 (3%)	35	62

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	36	GLU
1	A	61	ARG
1	A	78	GLN
1	A	138	ASP
1	A	144	GLU

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Mol	Chain	Res	Type
1	A	167	GLN
1	A	245	VAL
1	A	306	LEU
1	B	36	GLU
1	B	79	SER
1	B	138	ASP
1	B	162	GLN
1	B	237	ASN
1	B	291	GLN
1	B	307	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	119	HIS
1	A	213	HIS
1	A	276	GLN
1	B	60	GLN
1	B	119	HIS
1	B	260	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](#)

Of 10 ligands modelled in this entry, 3 are monoatomic - leaving 7 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and

the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	GOL	A	404	-	5,5,5	0.38	0	5,5,5	0.29	0
2	NAG	A	403	1	14,14,15	0.39	0	17,19,21	0.33	0
2	NAG	B	402	1	14,14,15	0.21	0	17,19,21	0.45	0
2	NAG	B	401	1	14,14,15	0.39	0	17,19,21	0.65	1 (5%)
2	NAG	A	401	1	14,14,15	0.19	0	17,19,21	0.56	0
2	NAG	A	402	1	14,14,15	0.28	0	17,19,21	0.43	0
3	GOL	B	403	-	5,5,5	0.38	0	5,5,5	0.32	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	GOL	A	404	-	-	2/4/4/4	-
2	NAG	A	403	1	-	2/6/23/26	0/1/1/1
2	NAG	B	402	1	-	0/6/23/26	0/1/1/1
2	NAG	B	401	1	-	0/6/23/26	0/1/1/1
2	NAG	A	401	1	-	2/6/23/26	0/1/1/1
2	NAG	A	402	1	-	0/6/23/26	0/1/1/1
3	GOL	B	403	-	-	2/4/4/4	-

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	401	NAG	C1-O5-C5	2.26	115.21	112.19

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	404	GOL	O1-C1-C2-C3
2	A	401	NAG	O5-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
2	A	401	NAG	C4-C5-C6-O6
3	B	403	GOL	O1-C1-C2-C3
3	A	404	GOL	O1-C1-C2-O2
3	B	403	GOL	O1-C1-C2-O2
2	A	403	NAG	C1-C2-N2-C7
2	A	403	NAG	C4-C5-C6-O6

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	402	NAG	1	0
2	B	401	NAG	1	0

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	289/334 (86%)	0.86	29 (10%)	12	10	21, 44, 78, 113	0
1	B	284/334 (85%)	1.04	32 (11%)	10	8	26, 52, 83, 116	0
All	All	573/668 (85%)	0.95	61 (10%)	11	9	21, 48, 82, 116	0

All (61) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	294	ALA	5.0
1	A	32	VAL	4.4
1	B	289	LEU	4.4
1	A	310	ALA	4.2
1	A	309	CYS	4.0
1	B	308	GLY	3.8
1	A	27	CYS	3.6
1	B	293	LEU	3.5
1	B	202	HIS	3.5
1	A	333	PRO	3.4
1	B	335	CYS	3.4
1	B	30	ALA	3.4
1	B	75	ALA	3.4
1	B	29	GLY	3.3
1	B	54	GLY	3.3
1	A	163	ASP	3.3
1	B	179	ASN	3.2
1	B	71	HIS	3.2
1	B	288	ASN	3.2
1	B	302	ALA	3.1
1	A	79	SER	3.1
1	A	169	LEU	3.1
1	B	263	VAL	3.0
1	A	78	GLN	3.0

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Mol	Chain	Res	Type	RSRZ
1	A	179	ASN	3.0
1	A	303	ALA	3.0
1	B	239	SER	2.9
1	A	30	ALA	2.9
1	A	71	HIS	2.8
1	A	33	CYS	2.7
1	B	163	ASP	2.6
1	A	26	PRO	2.6
1	A	81	ARG	2.6
1	B	26	PRO	2.6
1	B	81	ARG	2.5
1	B	259	ASP	2.5
1	B	295	ASP	2.5
1	B	306	LEU	2.5
1	B	82	ASN	2.5
1	A	28	PRO	2.4
1	A	335	CYS	2.4
1	A	307	GLU	2.4
1	A	102	THR	2.3
1	B	36	GLU	2.3
1	B	267	ARG	2.3
1	B	37	PRO	2.3
1	A	176	ASP	2.2
1	B	307	GLU	2.2
1	A	302	ALA	2.2
1	A	289	LEU	2.2
1	A	237	ASN	2.2
1	A	53	THR	2.2
1	A	80	CYS	2.2
1	A	69	ILE	2.1
1	A	336	CYS	2.1
1	A	76	SER	2.1
1	B	189	ARG	2.1
1	B	304	SER	2.1
1	B	31	CYS	2.1
1	B	28	PRO	2.0
1	B	57	ALA	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 [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	NAG	A	403	14/15	0.39	0.26	84,98,104,105	0
2	NAG	B	402	14/15	0.48	0.21	72,82,87,87	0
2	NAG	A	402	14/15	0.69	0.18	64,77,86,86	0
2	NAG	B	401	14/15	0.70	0.17	56,64,72,73	0
2	NAG	A	401	14/15	0.74	0.15	46,56,61,62	0
3	GOL	B	403	6/6	0.76	0.15	43,50,53,55	0
3	GOL	A	404	6/6	0.78	0.18	59,63,65,67	0
4	CL	B	404	1/1	0.85	0.18	77,77,77,77	0
4	CL	A	405	1/1	0.94	0.14	67,67,67,67	0
4	CL	B	405	1/1	0.97	0.15	61,61,61,61	0

6.5 Other polymers [i](#)

There are no such residues in this entry.