



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 01:33 PM GMT

PDB ID : 3U2P  
Title : Crystal structure of N-terminal three extracellular domains of ErbB4/Her4  
Authors : Liu, P.; Bouyain, S.; Elgenbrot, C.; Leahy, D.J.  
Deposited on : 2011-10-04  
Resolution : 2.57 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](mailto:validation@mail.wwpdb.org)  
A user guide is available at  
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>  
with specific help available everywhere you see the ⓘ symbol.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
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) : trunk26865

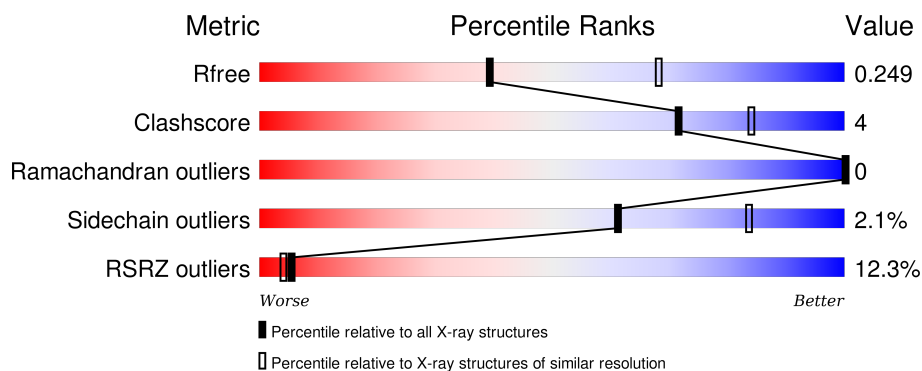
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*


The reported resolution of this entry is 2.57 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	2636 (2.60-2.56)
Clashscore	102246	3003 (2.60-2.56)
Ramachandran outliers	100387	2956 (2.60-2.56)
Sidechain outliers	100360	2956 (2.60-2.56)
RSRZ outliers	91569	2642 (2.60-2.56)

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  $\geq 3$ , 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  $\leq 5\%$ . 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	497	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	NAG	A	1002	-	-	-	X

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	NAG	A	1004	-	-	-	X

## 2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 3893 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 Receptor tyrosine-protein kinase erbB-4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	479	Total	C	N	O	S	0	0	0
			3737	2331	650	715	41			

- Molecule 2 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: 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	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		

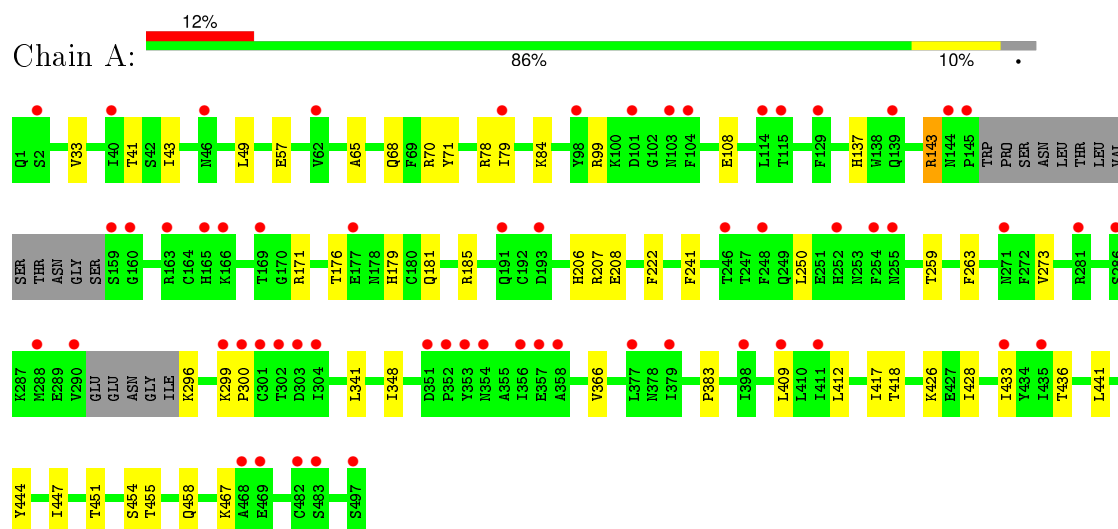
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	86	Total	O	0	0
			86	86		

### 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: Receptor tyrosine-protein kinase erbB-4



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	42.17Å 85.56Å 152.40Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	24.87 – 2.57 24.87 – 2.57	Depositor EDS
% Data completeness (in resolution range)	(Not available) (24.87-2.57) 96.9 (24.87-2.57)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.91 (at 2.57Å)	Xtriage
Refinement program	BUSTER 2.8.0	Depositor
R, $R_{free}$	0.192 , 0.243 0.196 , 0.249	Depositor DCC
$R_{free}$ test set	920 reflections (5.46%)	DCC
Wilson B-factor (Å <sup>2</sup> )	44.6	Xtriage
Anisotropy	0.058	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 56.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 17764 reflections	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	3893	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.45% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

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

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	0.50	0/3811	0.75	0/5160

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 [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	3737	0	3589	32	0
2	A	70	0	64	1	0
3	A	86	0	0	0	0
All	All	3893	0	3653	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:137:HIS:H	1:A:181:GLN:HE22	1.24	0.83

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:176:THR:H	1:A:179:HIS:HD2	1.34	0.72
1:A:409:LEU:HB3	1:A:433:ILE:HD12	1.76	0.67
1:A:78:ARG:HD3	1:A:222:PHE:HB3	1.80	0.64
1:A:68:GLN:HG2	1:A:99:ARG:HB2	1.82	0.61
1:A:454:SER:H	1:A:458:GLN:HE22	1.51	0.59
1:A:206:HIS:HD2	1:A:208:GLU:H	1.52	0.56
1:A:43:ILE:HG21	1:A:49:LEU:HD11	1.87	0.56
1:A:455:THR:H	1:A:458:GLN:HE21	1.52	0.56
1:A:428:ILE:HD13	1:A:433:ILE:HD11	1.87	0.56
1:A:273:VAL:HG12	1:A:296:LYS:HB3	1.89	0.54
1:A:417:ILE:HD12	1:A:441:LEU:HD13	1.91	0.53
1:A:176:THR:H	1:A:179:HIS:CD2	2.22	0.52
1:A:143:ARG:HA	1:A:185:ARG:HH12	1.75	0.52
1:A:467:LYS:NZ	2:A:1005:NAG:H82	2.24	0.51
1:A:71:TYR:HD1	1:A:108:GLU:HG3	1.75	0.50
1:A:241:PHE:HB3	1:A:250:LEU:HB3	1.96	0.47
1:A:33:VAL:HG21	1:A:263:PHE:HZ	1.81	0.46
1:A:455:THR:H	1:A:458:GLN:NE2	2.12	0.46
1:A:299:LYS:HD2	1:A:300:PRO:HD2	1.98	0.46
1:A:137:HIS:HB2	1:A:171:ARG:HD3	1.97	0.45
1:A:57:GLU:HG3	1:A:79:ILE:HG23	1.99	0.44
1:A:143:ARG:HA	1:A:185:ARG:NH1	2.33	0.44
1:A:348:ILE:HD12	1:A:383:PRO:HD3	1.99	0.43
1:A:412:LEU:HD12	1:A:436:THR:HG22	2.01	0.43
1:A:206:HIS:CD2	1:A:208:GLU:H	2.35	0.43
1:A:41:THR:HA	1:A:65:ALA:O	2.19	0.42
1:A:341:LEU:HA	1:A:341:LEU:HD23	1.88	0.42
1:A:444:TYR:HA	1:A:447:ILE:HD12	2.00	0.42
1:A:71:TYR:CD1	1:A:108:GLU:HG3	2.55	0.42
1:A:454:SER:N	1:A:458:GLN:HE22	2.18	0.42
1:A:428:ILE:HG21	1:A:433:ILE:HD11	2.03	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	473/497 (95%)	449 (95%)	24 (5%)	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	423/439 (96%)	414 (98%)	9 (2%)	61	84

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

Mol	Chain	Res	Type
1	A	70	ARG
1	A	84	LYS
1	A	143	ARG
1	A	207	ARG
1	A	259	THR
1	A	366	VAL
1	A	418	THR
1	A	426	LYS
1	A	451	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	179	HIS
1	A	181	GLN
1	A	206	HIS
1	A	421	GLN
1	A	440	ASN
1	A	457	ASN
1	A	458	GLN

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Mol	Chain	Res	Type
1	A	479	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 carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

5 ligands are modelled in this entry.

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$
2	NAG	A	1001	1	14,14,15	1.60	2 (14%)	15,19,21	3.60	7 (46%)
2	NAG	A	1002	1	14,14,15	1.26	1 (7%)	15,19,21	1.26	1 (6%)
2	NAG	A	1003	1	14,14,15	1.60	3 (21%)	15,19,21	1.29	1 (6%)
2	NAG	A	1004	1	14,14,15	2.12	4 (28%)	15,19,21	2.16	5 (33%)
2	NAG	A	1005	1	14,14,15	1.56	2 (14%)	15,19,21	1.86	4 (26%)

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
2	NAG	A	1001	1	-	0/6/23/26	0/1/1/1
2	NAG	A	1002	1	-	0/6/23/26	0/1/1/1
2	NAG	A	1003	1	-	0/6/23/26	0/1/1/1
2	NAG	A	1004	1	-	0/6/23/26	0/1/1/1
2	NAG	A	1005	1	-	0/6/23/26	0/1/1/1

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1001	NAG	C2-N2	-2.18	1.42	1.46
2	A	1004	NAG	C4-C5	2.07	1.57	1.53
2	A	1003	NAG	C2-N2	2.16	1.50	1.46
2	A	1003	NAG	C3-C2	2.17	1.57	1.52
2	A	1004	NAG	C2-N2	2.29	1.50	1.46
2	A	1005	NAG	C1-C2	2.41	1.55	1.52
2	A	1002	NAG	C1-C2	2.78	1.56	1.52
2	A	1005	NAG	C4-C3	2.86	1.59	1.52
2	A	1004	NAG	C3-C2	2.96	1.59	1.52
2	A	1003	NAG	C1-C2	4.40	1.58	1.52
2	A	1001	NAG	C1-C2	4.66	1.58	1.52
2	A	1004	NAG	C1-C2	5.38	1.59	1.52

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1001	NAG	C3-C2-N2	-4.96	98.68	110.56
2	A	1001	NAG	C8-C7-N2	-3.99	108.47	116.11
2	A	1004	NAG	C3-C2-N2	-2.95	103.48	110.56
2	A	1004	NAG	O4-C4-C3	-2.12	105.55	110.34
2	A	1001	NAG	O4-C4-C5	-2.02	103.88	109.24
2	A	1002	NAG	O6-C6-C5	2.18	118.53	111.33
2	A	1005	NAG	O3-C3-C4	2.19	115.27	110.34
2	A	1005	NAG	O6-C6-C5	2.48	119.52	111.33
2	A	1004	NAG	O3-C3-C2	3.09	115.23	109.11
2	A	1005	NAG	O5-C5-C6	3.09	114.03	107.35
2	A	1003	NAG	C4-C3-C2	3.19	116.18	111.23
2	A	1004	NAG	C1-O5-C5	3.34	116.48	112.25
2	A	1001	NAG	O7-C7-C8	3.45	128.39	122.06
2	A	1005	NAG	C1-O5-C5	3.76	117.03	112.25
2	A	1001	NAG	C4-C3-C2	4.35	118.00	111.23
2	A	1001	NAG	C1-O5-C5	4.77	118.30	112.25
2	A	1004	NAG	C4-C3-C2	5.57	119.89	111.23
2	A	1001	NAG	C2-N2-C7	9.49	135.24	123.04

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1005	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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	479/497 (96%)	0.55	59 (12%) 5 4	24, 47, 82, 114	0

All (59) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	159	SER	9.0
1	A	353	TYR	7.1
1	A	352	PRO	5.0
1	A	303	ASP	4.8
1	A	160	GLY	4.6
1	A	101	ASP	4.2
1	A	145	PRO	4.0
1	A	144	ASN	4.0
1	A	483	SER	3.8
1	A	299	LYS	3.8
1	A	286	SER	3.7
1	A	252	HIS	3.7
1	A	377	LEU	3.7
1	A	302	THR	3.7
1	A	254	PHE	3.6
1	A	103	ASN	3.6
1	A	356	ILE	3.5
1	A	290	VAL	3.4
1	A	191	GLN	3.4
1	A	358	ALA	3.4
1	A	248	PHE	3.2
1	A	271	ASN	3.1
1	A	281	ARG	3.1
1	A	177	GLU	3.0
1	A	357	GLU	3.0
1	A	288	MET	2.8
1	A	409	LEU	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	40	ILE	2.7
1	A	62	VAL	2.7
1	A	46	ASN	2.6
1	A	139	GLN	2.6
1	A	166	LYS	2.5
1	A	163	ARG	2.5
1	A	246	THR	2.5
1	A	497	SER	2.5
1	A	354	ASN	2.4
1	A	169	THR	2.4
1	A	79	ILE	2.4
1	A	129	PHE	2.4
1	A	469	GLU	2.4
1	A	304	ILE	2.4
1	A	114	LEU	2.4
1	A	398	ILE	2.3
1	A	435	ILE	2.3
1	A	255	ASN	2.3
1	A	379	ILE	2.2
1	A	115	THR	2.2
1	A	351	ASP	2.2
1	A	411	ILE	2.2
1	A	193	ASP	2.2
1	A	301	CYS	2.1
1	A	482	CYS	2.1
1	A	468	ALA	2.1
1	A	98	TYR	2.1
1	A	104	PHE	2.1
1	A	2	SER	2.0
1	A	165	HIS	2.0
1	A	433	ILE	2.0
1	A	300	PRO	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 [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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> 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	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
2	NAG	A	1004	14/15	0.66	0.34	7.44	90,94,97,97	0
2	NAG	A	1002	14/15	0.83	0.31	3.18	58,61,64,65	0
2	NAG	A	1001	14/15	0.96	0.14	-	43,47,50,51	0
2	NAG	A	1005	14/15	0.68	0.34	-	94,98,100,100	0
2	NAG	A	1003	14/15	0.89	0.23	-	65,68,71,72	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.