



Full wwPDB X-ray Structure Validation Report ⓘ

May 23, 2016 – 09:52 PM EDT

PDB ID : 4RT6
Title : Structure of a complex between hemopexin and hemopexin binding protein
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Deposited on : 2014-11-13
Resolution : 2.80 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
<http://wwpdb.org/validation/2016/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
Mogul : 1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20027674
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) : rb-20027674

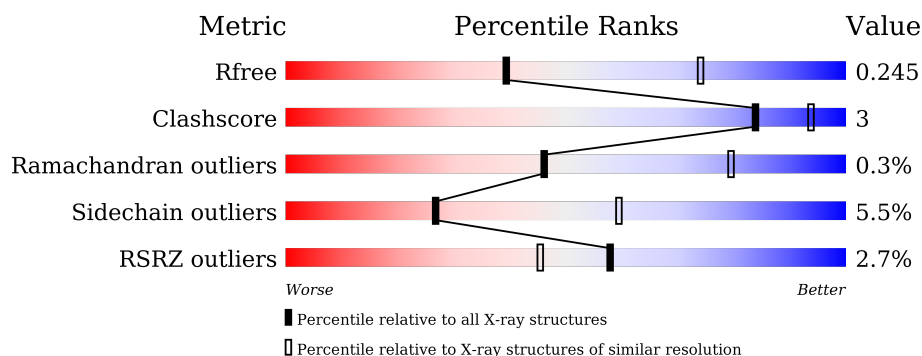
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	2393 (2.80-2.80)
Clashscore	102246	2827 (2.80-2.80)
Ramachandran outliers	100387	2782 (2.80-2.80)
Sidechain outliers	100360	2784 (2.80-2.80)
RSRZ outliers	91569	2404 (2.80-2.80)

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. 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	884	<div> <div>2%</div> <div> <div></div> <div>83%</div> <div>7%</div> <div>•</div> <div>9%</div> </div> </div>
2	B	214	<div> <div>6%</div> <div> <div></div> <div>57%</div> <div>21%</div> <div>•</div> <div>20%</div> </div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7841 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 Heme/hemopexin-binding protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	805	Total	C	N	O	S	0	0	0
			6204	3833	1107	1252	12			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	855	SER	CYS	engineered mutation	UNP P44602
A	861	SER	CYS	engineered mutation	UNP P44602

- Molecule 2 is a protein called Hemopexin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	172	Total	C	N	O	S	0	0	0
			1417	918	237	255	7			

- Molecule 3 is a polymer of unknown type called SUGAR (2-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	B	2	Total	C	N	O	0	0
			28	16	2	10		

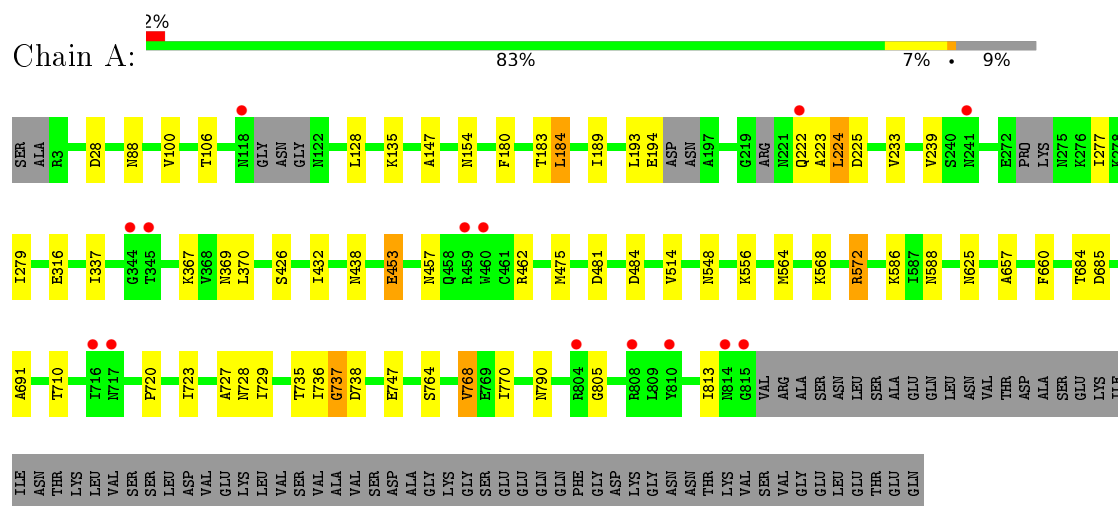
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	174	Total	O	0	0
			174	174		
4	B	18	Total	O	0	0
			18	18		

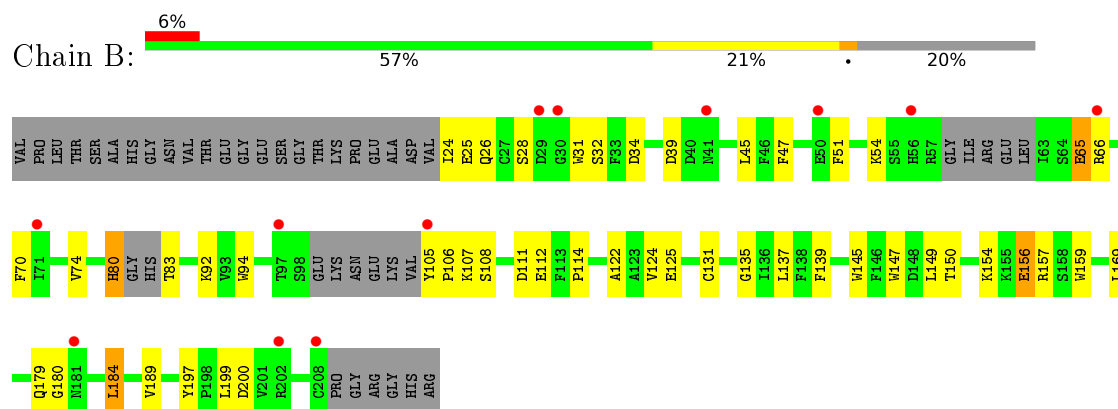
3 Residue-property plots [i](#)

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: Heme/hemopexin-binding protein



• Molecule 2: Hemopexin



4 Data and refinement statistics

Property	Value	Source
Space group	P 2 ₁ 2 ₁ 2	Depositor
Cell constants a, b, c, α , β , γ	45.06 Å 348.86 Å 77.35 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	24.43 – 2.80 24.43 – 2.80	Depositor EDS
% Data completeness (in resolution range)	96.0 (24.43-2.80) 96.3 (24.43-2.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.63 (at 2.80 Å)	Xtriage
Refinement program	BUSTER 2.10.1	Depositor
R, R_{free}	0.180 , 0.230 0.195 , 0.245	Depositor DCC
R_{free} test set	1517 reflections (5.33%)	DCC
Wilson B-factor (Å ²)	51.7	Xtriage
Anisotropy	0.095	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 55.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	7841	wwPDB-VP
Average B, all atoms (Å ²)	55.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.43% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: 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	0.51	0/6276	0.73	0/8448
2	B	0.48	0/1464	0.69	0/1984
All	All	0.50	0/7740	0.73	0/10432

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	6204	0	6167	27	0
2	B	1417	0	1317	20	0
3	B	28	0	25	0	0
4	A	174	0	0	4	0
4	B	18	0	0	0	0
All	All	7841	0	7509	46	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 (46) 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:239:VAL:HG22	4:A:1061:HOH:O	1.79	0.80
2:B:28:SER:HB2	2:B:31:TRP:HB2	1.79	0.64
1:A:768:VAL:O	1:A:768:VAL:HG12	1.99	0.63
2:B:80:HIS:CE1	2:B:125:GLU:HB2	2.35	0.61
2:B:92:LYS:HB2	2:B:94:TRP:CZ3	2.41	0.56
2:B:114:PRO:HD2	2:B:147:TRP:HH2	1.70	0.55
2:B:114:PRO:HD2	2:B:147:TRP:CH2	2.43	0.54
2:B:122:ALA:HB3	2:B:139:PHE:HB2	1.91	0.53
2:B:184:LEU:HA	2:B:197:TYR:CD1	2.45	0.52
1:A:514:VAL:CG1	1:A:572:ARG:HG3	2.40	0.52
1:A:438:ASN:HA	1:A:481:ASP:O	2.10	0.52
1:A:805:GLY:HA3	4:A:1047:HOH:O	2.10	0.51
1:A:100:VAL:O	1:A:147:ALA:HA	2.11	0.51
2:B:34:ASP:HB2	2:B:47:PHE:O	2.10	0.50
1:A:586:LYS:HE2	1:A:588:ASN:HB2	1.92	0.50
1:A:183:THR:HG22	1:A:184:LEU:H	1.77	0.50
1:A:128:LEU:HD11	1:A:194:GLU:HB2	1.95	0.49
2:B:108:SER:HB3	2:B:111:ASP:HB2	1.95	0.49
2:B:34:ASP:HB3	2:B:74:VAL:O	2.13	0.48
1:A:180:PHE:CZ	1:A:189:ILE:HD12	2.50	0.47
1:A:685:ASP:OD2	1:A:737:GLY:HA3	2.13	0.47
1:A:106:THR:HA	1:A:154:ASN:O	2.15	0.47
1:A:691:ALA:HA	1:A:747:GLU:O	2.15	0.46
2:B:32:SER:HA	2:B:180:GLY:HA2	1.96	0.46
1:A:727:ALA:O	1:A:728:ASN:HB2	2.16	0.46
1:A:432:ILE:O	1:A:475:MET:HA	2.16	0.45
1:A:367:LYS:HE2	1:A:369:ASN:OD1	2.16	0.45
1:A:337:ILE:HD11	1:A:370:LEU:HD21	1.98	0.45
1:A:453:GLU:H	1:A:453:GLU:HG3	1.56	0.45
1:A:736:ILE:CD1	1:A:770:ILE:HG12	2.47	0.44
2:B:145:TRP:NE1	2:B:156:GLU:HG3	2.33	0.43
1:A:88:ASN:HA	1:A:135:LYS:HE2	2.01	0.43
1:A:764:SER:HB2	1:A:770:ILE:HG13	2.01	0.43
2:B:135:GLY:HA2	2:B:149:LEU:HG	2.01	0.42
1:A:223:ALA:O	1:A:224:LEU:HB2	2.19	0.42
2:B:157:ARG:HH11	2:B:159:TRP:HZ2	1.67	0.42
2:B:65:GLU:HG3	2:B:65:GLU:H	1.59	0.42
2:B:94:TRP:CE2	2:B:106:PRO:HB3	2.55	0.41
2:B:70:PHE:HE2	2:B:74:VAL:HG22	1.85	0.41
1:A:710:THR:HG23	1:A:729:ILE:HA	2.02	0.41
1:A:233:VAL:HG23	4:A:1039:HOH:O	2.21	0.41
2:B:24:ILE:HG23	2:B:25:GLU:H	1.86	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:720:PRO:HB2	2:B:197:TYR:OH	2.21	0.40
1:A:625:ASN:HA	1:A:657:ALA:O	2.21	0.40
2:B:39:ASP:HB3	2:B:45:LEU:HD11	2.04	0.40
1:A:548:ASN:HB2	4:A:974:HOH:O	2.21	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	A	795/884 (90%)	759 (96%)	33 (4%)	3 (0%)	39	74
2	B	164/214 (77%)	149 (91%)	15 (9%)	0	100	100
All	All	959/1098 (87%)	908 (95%)	48 (5%)	3 (0%)	46	79

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	768	VAL
1	A	737	GLY
1	A	224	LEU

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	690/756 (91%)	666 (96%)	24 (4%)	43	77
2	B	151/184 (82%)	129 (85%)	22 (15%)	4	11
All	All	841/940 (90%)	795 (94%)	46 (6%)	27	59

All (46) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	28	ASP
1	A	184	LEU
1	A	193	LEU
1	A	222	GLN
1	A	225	ASP
1	A	277	ILE
1	A	279	ILE
1	A	316	GLU
1	A	426	SER
1	A	453	GLU
1	A	457	ASN
1	A	462	ARG
1	A	484	ASP
1	A	556	LYS
1	A	564	MET
1	A	568	LYS
1	A	572	ARG
1	A	660	PHE
1	A	684	THR
1	A	723	ILE
1	A	735	THR
1	A	738	ASP
1	A	790	ASN
1	A	813	ILE
2	B	26	GLN
2	B	51	PHE
2	B	54	LYS
2	B	65	GLU
2	B	66	ARG
2	B	80	HIS
2	B	83	THR
2	B	105	TYR
2	B	107	LYS
2	B	112	GLU
2	B	124	VAL

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Mol	Chain	Res	Type
2	B	131	CYS
2	B	137	LEU
2	B	150	THR
2	B	154	LYS
2	B	156	GLU
2	B	169	LEU
2	B	179	GLN
2	B	184	LEU
2	B	189	VAL
2	B	199	LEU
2	B	200	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	73	GLN
1	A	122	ASN
1	A	256	ASN
1	A	282	GLN
1	A	581	ASN
1	A	791	GLN
2	B	41	ASN

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 ⓘ

2 carbohydrates 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	$\# Z > 2$
3	NAG	B	301	3,2	14,14,15	0.28	0	15,19,21	0.86	1 (6%)
3	NAG	B	302	3	14,14,15	0.31	0	15,19,21	0.69	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	NAG	B	301	3,2	-	0/6/23/26	0/1/1/1
3	NAG	B	302	3	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	301	NAG	C1-O5-C5	3.10	116.70	112.14

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	805/884 (91%)	-0.34	14 (1%) 73 63	25, 46, 77, 121	1 (0%)
2	B	172/214 (80%)	0.11	12 (6%) 19 11	45, 75, 115, 129	0
All	All	977/1098 (88%)	-0.26	26 (2%) 58 45	25, 50, 95, 129	1 (0%)

All (26) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	815	GLY	4.8
2	B	105	TYR	4.2
1	A	344	GLY	3.8
2	B	181	ASN	3.6
1	A	345	THR	3.4
2	B	71	ILE	3.4
1	A	460	TRP	3.4
2	B	30	GLY	3.2
1	A	814	ASN	3.0
2	B	66	ARG	2.8
2	B	56	HIS	2.6
1	A	716	ILE	2.6
2	B	41	ASN	2.5
1	A	241	ASN	2.5
2	B	97	THR	2.5
1	A	808	ARG	2.4
1	A	118	ASN	2.4
2	B	202	ARG	2.3
1	A	459	ARG	2.2
1	A	222	GLN	2.2
2	B	29	ASP	2.2
2	B	50	GLU	2.1
1	A	810	TYR	2.1
2	B	208	CYS	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	804	ARG	2.1
1	A	717	ASN	2.1

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](#)

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, 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	LLDF	B-factors(Å ²)	Q<0.9
3	NAG	B	302	14/15	0.70	0.51	-	118,123,129,130	0
3	NAG	B	301	14/15	0.85	0.43	-	110,114,120,122	0

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.