



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 01:20 PM UTC

PDB ID : 6GRF / pdb_00006grf
Title : Crystal structure of the tandem DUF26 ectodomain from the Arabidopsis thaliana cysteine-rich receptor-like protein PDLP8.
Authors : Brandt, B.; Hothorn, M.
Deposited on : 2018-06-11
Resolution : 1.95 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

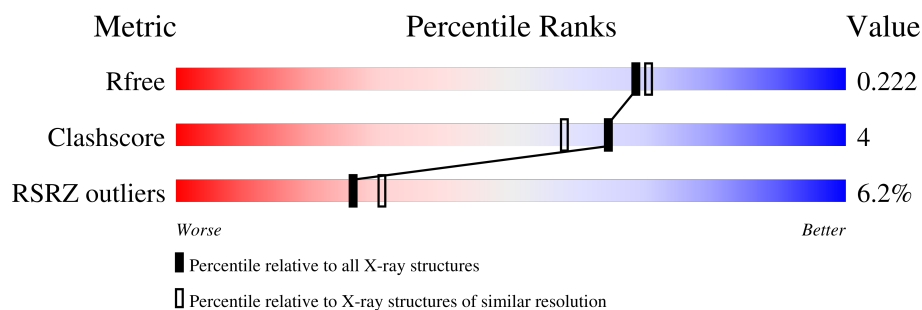
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.95 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	3494 (1.96-1.96)
Clashscore	190562	3612 (1.96-1.96)
RSRZ outliers	180081	3495 (1.96-1.96)

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	242	 6% 76% 11% 13%
1	B	242	 6% 82% 5% 13%
1	C	242	 4% 76% 9% 15%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5042 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 Cysteine-rich repeat secretory protein 15.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	210	Total	C	N	O	S	0	4	0
			1624	1012	265	335	12			
1	B	211	Total	C	N	O	S	0	1	0
			1611	998	265	336	12			
1	C	206	Total	C	N	O	S	0	1	0
			1586	986	263	325	12			

There are 30 discrepancies between the modelled and reference sequences:

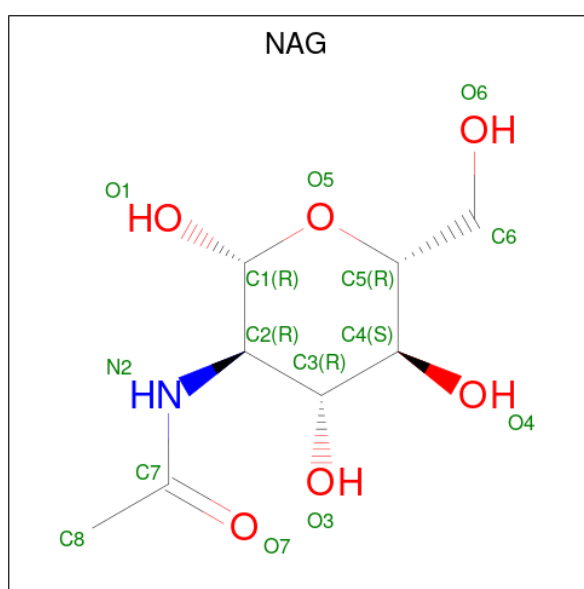
Chain	Residue	Modelled	Actual	Comment	Reference
A	254	LEU	-	expression tag	UNP Q6NKQ9
A	255	GLU	-	expression tag	UNP Q6NKQ9
A	256	GLY	-	expression tag	UNP Q6NKQ9
A	257	SER	-	expression tag	UNP Q6NKQ9
A	258	GLU	-	expression tag	UNP Q6NKQ9
A	259	ASN	-	expression tag	UNP Q6NKQ9
A	260	LEU	-	expression tag	UNP Q6NKQ9
A	261	TYR	-	expression tag	UNP Q6NKQ9
A	262	PHE	-	expression tag	UNP Q6NKQ9
A	263	GLN	-	expression tag	UNP Q6NKQ9
B	254	LEU	-	expression tag	UNP Q6NKQ9
B	255	GLU	-	expression tag	UNP Q6NKQ9
B	256	GLY	-	expression tag	UNP Q6NKQ9
B	257	SER	-	expression tag	UNP Q6NKQ9
B	258	GLU	-	expression tag	UNP Q6NKQ9
B	259	ASN	-	expression tag	UNP Q6NKQ9
B	260	LEU	-	expression tag	UNP Q6NKQ9
B	261	TYR	-	expression tag	UNP Q6NKQ9
B	262	PHE	-	expression tag	UNP Q6NKQ9
B	263	GLN	-	expression tag	UNP Q6NKQ9
C	254	LEU	-	expression tag	UNP Q6NKQ9
C	255	GLU	-	expression tag	UNP Q6NKQ9
C	256	GLY	-	expression tag	UNP Q6NKQ9

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Chain	Residue	Modelled	Actual	Comment	Reference
C	257	SER	-	expression tag	UNP Q6NKKQ9
C	258	GLU	-	expression tag	UNP Q6NKKQ9
C	259	ASN	-	expression tag	UNP Q6NKKQ9
C	260	LEU	-	expression tag	UNP Q6NKKQ9
C	261	TYR	-	expression tag	UNP Q6NKKQ9
C	262	PHE	-	expression tag	UNP Q6NKKQ9
C	263	GLN	-	expression tag	UNP Q6NKKQ9

- 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	B	1	Total	C	N	O	0	0
			14	8	1	5		

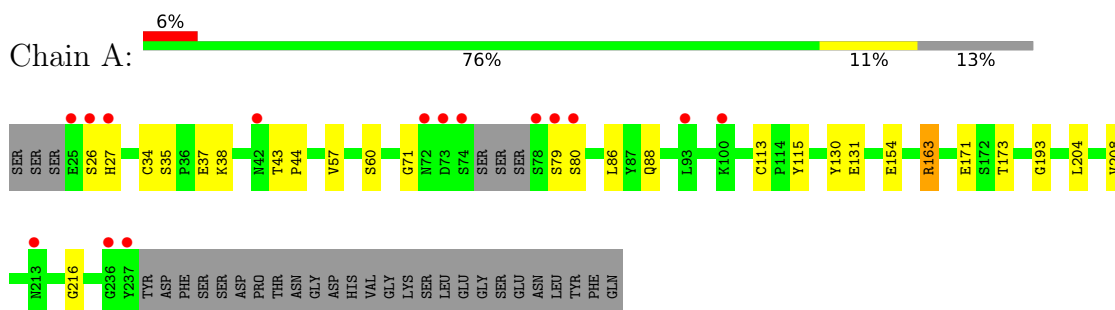
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	74	Total	O	0	0
			74	74		
3	B	64	Total	O	0	0
			64	64		
3	C	55	Total	O	0	0
			55	55		

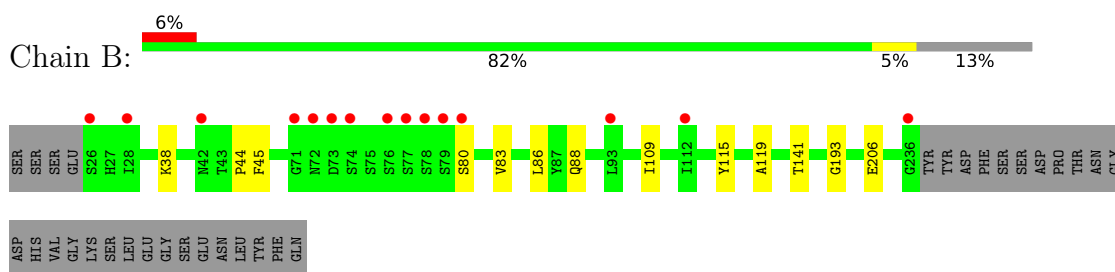
3 Residue-property plots [i](#)

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.

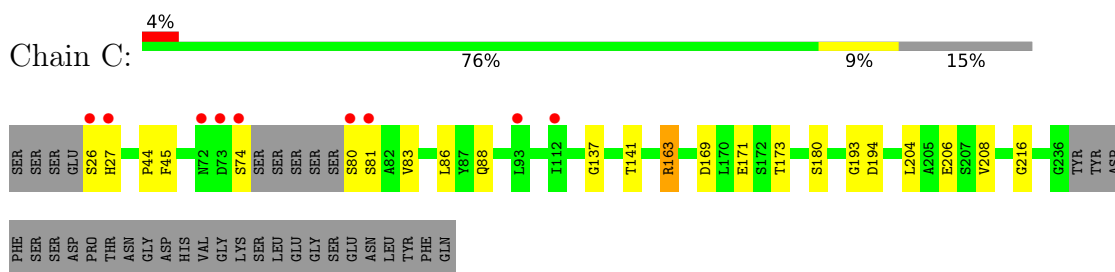
- Molecule 1: Cysteine-rich repeat secretory protein 15



- Molecule 1: Cysteine-rich repeat secretory protein 15



- Molecule 1: Cysteine-rich repeat secretory protein 15



4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	143.86Å 143.86Å 59.72Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.10 – 1.95 20.10 – 1.95	Depositor EDS
% Data completeness (in resolution range)	99.8 (20.10-1.95) 99.8 (20.10-1.95)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.31	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.02 (at 1.94Å)	Xtriage
Refinement program	REFMAC 5.8.0222	Depositor
R, R_{free}	0.229 , 0.262 (Not available) , 0.222	Depositor DCC
R_{free} test set	2591 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	30.8	Xtriage
Anisotropy	0.488	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 28.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.000 for -h,-k,l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	5042	wwPDB-VP
Average B, all atoms (Å ²)	43.0	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 55.73 % 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 3.0343e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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

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	0.73	0/1667	0.97	0/2252
1	B	0.67	0/1645	0.97	0/2224
1	C	0.69	0/1619	0.98	0/2187
All	All	0.70	0/4931	0.97	0/6663

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	2
1	C	0	1
All	All	0	3

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	163	ARG	Sidechain
1	A	71	GLY	Peptide
1	C	163	ARG	Sidechain

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	1624	0	1529	15	1
1	B	1611	0	1510	7	0
1	C	1586	0	1493	14	0
2	A	14	0	13	0	0
2	B	14	0	13	0	0
3	A	74	0	0	0	0
3	B	64	0	0	0	0
3	C	55	0	0	1	0
All	All	5042	0	4558	36	1

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 (36) 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:34:CYS:HG	1:A:113:CYS:HG	1.09	0.92
1:A:34:CYS:CB	1:A:113:CYS:HG	2.08	0.67
1:A:88:GLN:HE21	1:A:193:GLY:H	1.49	0.60
1:C:86:LEU:C	1:C:86:LEU:HD12	2.32	0.55
1:B:109:ILE:HD11	1:B:119:ALA:HB1	1.88	0.54
1:A:130:TYR:C	1:A:131:GLU:HG2	2.31	0.54
1:C:88:GLN:HE21	1:C:193:GLY:H	1.56	0.54
1:A:44:PRO:HB2	1:A:80:SER:HA	1.91	0.53
1:C:171:GLU:HG3	1:C:208:VAL:HG21	1.92	0.50
1:C:44:PRO:HB2	1:C:80:SER:HA	1.93	0.49
1:C:26:SER:HA	3:C:349:HOH:O	2.12	0.49
1:B:44:PRO:HB2	1:B:80:SER:HA	1.96	0.48
1:C:137:GLY:HA2	1:C:194:ASP:O	2.16	0.46
1:B:45:PHE:CE2	1:B:83:VAL:HG22	2.51	0.46
1:A:26:SER:C	1:A:27:HIS:HD1	2.24	0.45
1:A:88:GLN:NE2	1:A:193:GLY:H	2.13	0.45
1:A:173:THR:HB	1:A:204:LEU:CD1	2.47	0.45
1:C:26:SER:C	1:C:27:HIS:HD1	2.23	0.44
1:C:74:SER:C	1:C:81:SER:OG	2.59	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:86:LEU:HD12	1:A:86:LEU:C	2.42	0.44
1:B:141:THR:HB	1:B:206:GLU:HG2	2.00	0.44
1:B:38:LYS:HA	1:B:115:TYR:O	2.18	0.42
1:A:171:GLU:HG3	1:A:208:VAL:HG21	2.01	0.42
1:B:88:GLN:HE21	1:B:193:GLY:H	1.67	0.42
1:A:57:VAL:O	1:A:60:SER:HB2	2.20	0.41
1:A:163:ARG:NH2	1:A:216:GLY:O	2.53	0.41
1:C:173:THR:HB	1:C:204:LEU:CD1	2.51	0.41
1:A:43:THR:CG2	1:A:79:SER:HB2	2.51	0.41
1:A:35:SER:HB2	1:A:37:GLU:OE1	2.20	0.41
1:A:38:LYS:HA	1:A:115:TYR:O	2.21	0.41
1:C:26:SER:C	1:C:27:HIS:ND1	2.79	0.40
1:C:163:ARG:NH2	1:C:216:GLY:O	2.54	0.40
1:C:141:THR:HB	1:C:206:GLU:HG2	2.03	0.40
1:C:169:ASP:OD2	1:C:180:SER:OG	2.37	0.40
1:B:86:LEU:C	1:B:86:LEU:HD12	2.47	0.40
1:C:45:PHE:CE2	1:C:83:VAL:HG22	2.57	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:154:GLU:OE2	1:A:154:GLU:OE2[4_554]	2.10	0.10

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

There are no protein backbone outliers to report in this entry.

5.3.2 Protein sidechains [i](#)

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

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](#)

2 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	B	301	1	14,14,15	0.48	0	17,19,21	0.82	0
2	NAG	A	301	1	14,14,15	0.40	0	17,19,21	0.92	1 (5%)

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	B	301	1	-	0/6/23/26	0/1/1/1
2	NAG	A	301	1	-	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(°)
2	A	301	NAG	C3-C4-C5	2.21	114.25	110.23

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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	210/242 (86%)	0.39	15 (7%)	22 25	20, 37, 68, 98	4 (1%)
1	B	211/242 (87%)	0.53	15 (7%)	22 25	26, 41, 69, 96	1 (0%)
1	C	206/242 (85%)	0.46	9 (4%)	39 45	23, 42, 66, 94	1 (0%)
All	All	627/726 (86%)	0.46	39 (6%)	26 31	20, 40, 69, 98	6 (0%)

All (39) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	26	SER	4.4
1	A	27	HIS	4.3
1	A	78	SER	4.1
1	C	93	LEU	3.8
1	A	236	GLY	3.7
1	A	42	ASN	3.7
1	A	79	SER	3.6
1	B	76	SER	3.6
1	B	78	SER	3.5
1	A	237	TYR	3.5
1	B	77	SER	3.2
1	B	236	GLY	3.2
1	A	73	ASP	3.1
1	A	26	SER	3.0
1	C	74	SER	3.0
1	C	72	ASN	3.0
1	A	74	SER	3.0
1	C	27	HIS	2.9
1	B	112	ILE	2.9
1	A	93	LEU	2.9
1	B	79	SER	2.8
1	C	80	SER	2.7
1	A	25	GLU	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	28	ILE	2.4
1	B	71	GLY	2.4
1	B	74	SER	2.4
1	B	80	SER	2.4
1	B	42	ASN	2.3
1	C	73	ASP	2.3
1	A	213	ASN	2.2
1	B	72	ASN	2.2
1	C	81	SER	2.2
1	B	73	ASP	2.1
1	A	80	SER	2.1
1	A	100[A]	LYS	2.1
1	C	112	ILE	2.1
1	B	26	SER	2.1
1	A	72	ASN	2.1
1	B	93	LEU	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](#)

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(Å ²)	Q<0.9
2	NAG	B	301	14/15	0.64	0.17	67,75,76,79	0
2	NAG	A	301	14/15	0.70	0.16	61,69,77,78	0

6.5 Other polymers [i](#)

There are no such residues in this entry.