



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 4, 2026 – 11:03 PM UTC

PDB ID : 3ZYG / pdb_00003zyg
Title : NETRING2 LAM AND EGF1 DOMAINS
Authors : Seiradake, E.; Coles, C.H.; Perestenko, P.V.; Harlos, K.; Mcilhinney, R.A.J.; Aricescu, A.R.; Jones, E.Y.
Deposited on : 2011-08-22
Resolution : 2.20 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

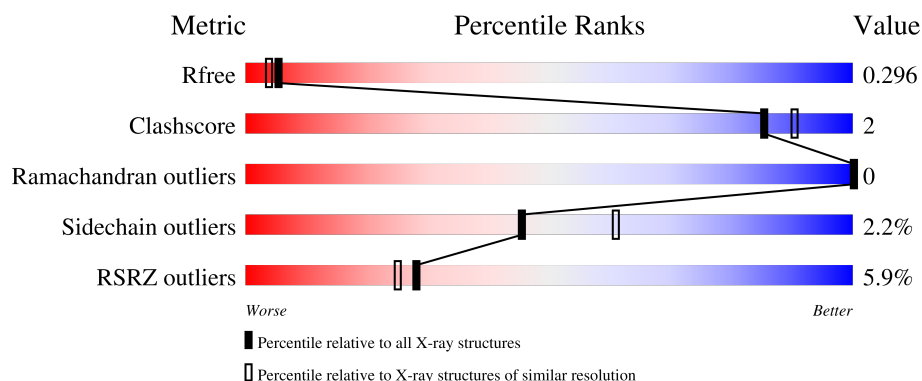
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	353	5% 84% 7% 9%
1	B	353	6% 80% 9% 10%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5383 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 NETRIN-G2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	322	Total	C	N	O	S	0	0	0
			2574	1619	451	478	26			
1	B	317	Total	C	N	O	S	0	0	0
			2543	1600	446	471	26			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	346	GLY	-	expression tag	UNP Q96CW9
A	347	THR	-	expression tag	UNP Q96CW9
A	348	HIS	-	expression tag	UNP Q96CW9
A	349	HIS	-	expression tag	UNP Q96CW9
A	350	HIS	-	expression tag	UNP Q96CW9
A	351	HIS	-	expression tag	UNP Q96CW9
A	352	HIS	-	expression tag	UNP Q96CW9
A	353	HIS	-	expression tag	UNP Q96CW9
B	346	GLY	-	expression tag	UNP Q96CW9
B	347	THR	-	expression tag	UNP Q96CW9
B	348	HIS	-	expression tag	UNP Q96CW9
B	349	HIS	-	expression tag	UNP Q96CW9
B	350	HIS	-	expression tag	UNP Q96CW9
B	351	HIS	-	expression tag	UNP Q96CW9
B	352	HIS	-	expression tag	UNP Q96CW9
B	353	HIS	-	expression tag	UNP Q96CW9

- Molecule 2 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆).



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	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 CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Ca	0	0
			1	1		
3	B	1	Total	Ca	0	0
			1	1		

- Molecule 4 is water.

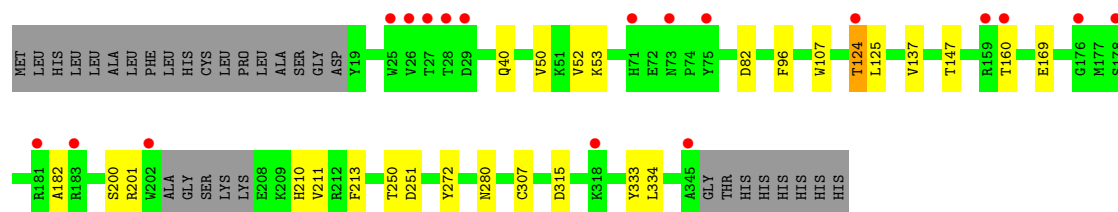
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	107	Total	O	0	0
			107	107		
4	B	101	Total	O	0	0
			101	101		

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.

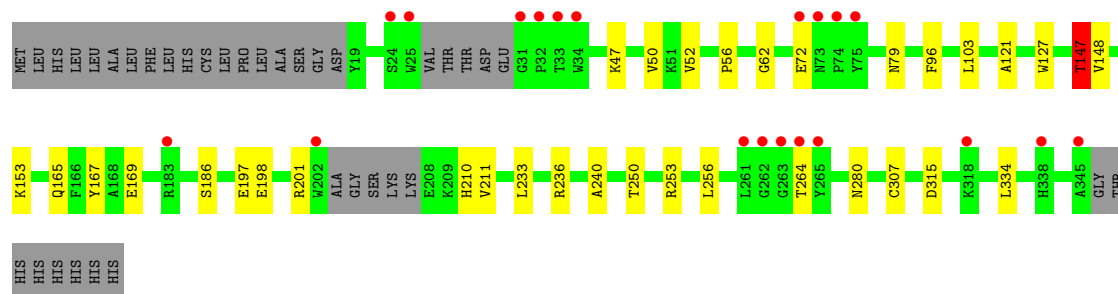
• Molecule 1: NETRIN-G2

Chain A: 



• Molecule 1: NETRIN-G2

Chain B: 



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	111.50Å 64.88Å 120.08Å 90.00° 103.06° 90.00°	Depositor
Resolution (Å)	30.00 – 2.20 30.00 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.1 (30.00-2.20) 99.1 (30.00-2.20)	Depositor EDS
R_{merge}	0.32	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.81 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.6.0077	Depositor
R, R_{free}	0.248 , 0.297 0.247 , 0.296	Depositor DCC
R_{free} test set	2125 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	18.9	Xtriage
Anisotropy	0.122	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 39.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	5383	wwPDB-VP
Average B, all atoms (Å ²)	25.0	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 55.23 % 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 3.2602e-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, CA

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	0.86	0/2643	1.19	16/3584 (0.4%)
1	B	0.82	0/2611	1.19	11/3537 (0.3%)
All	All	0.84	0/5254	1.19	27/7121 (0.4%)

There are no bond length outliers.

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	315	ASP	CA-CB-CG	6.75	119.35	112.60
1	A	210	HIS	N-CA-C	6.25	119.69	109.76
1	B	315	ASP	CA-CB-CG	6.16	118.76	112.60
1	A	307	CYS	N-CA-C	5.93	118.83	110.23
1	B	264	THR	N-CA-C	5.78	116.72	108.74
1	A	200	SER	CA-C-N	5.74	129.52	121.02
1	A	200	SER	C-N-CA	5.74	129.52	121.02
1	B	147	THR	N-CA-C	-5.74	105.38	112.90
1	B	211	VAL	N-CA-C	-5.73	100.14	108.17
1	A	147	THR	N-CA-C	-5.71	104.66	111.69
1	B	250	THR	N-CA-C	-5.71	106.86	113.88
1	A	250	THR	N-CA-C	-5.68	106.89	113.88
1	B	307	CYS	N-CA-C	5.61	118.47	110.10
1	A	124	THR	CB-CA-C	5.57	119.36	109.72
1	B	210	HIS	N-CA-C	5.55	119.29	109.96
1	A	182	ALA	CA-C-N	5.54	128.25	120.28
1	A	182	ALA	C-N-CA	5.54	128.25	120.28
1	B	62	GLY	N-CA-C	5.45	122.29	115.21
1	B	280	ASN	N-CA-C	5.37	117.24	108.55
1	A	107	TRP	N-CA-C	-5.36	102.55	110.48
1	A	211	VAL	N-CA-C	-5.33	100.70	108.17
1	A	280	ASN	N-CA-C	5.31	117.15	108.55

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	197	GLU	CA-C-N	5.27	128.08	120.38
1	B	197	GLU	C-N-CA	5.27	128.08	120.38
1	A	272	TYR	CA-C-N	5.13	129.40	120.68
1	A	272	TYR	C-N-CA	5.13	129.40	120.68
1	A	82	ASP	CA-CB-CG	5.05	117.66	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	A	2574	0	2427	10	0
1	B	2543	0	2399	12	0
2	A	28	0	26	0	0
2	B	28	0	26	3	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	107	0	0	0	0
4	B	101	0	0	0	0
All	All	5383	0	4878	22	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 2.

All (22) 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:52:VAL:HG11	1:A:96:PHE:CE2	1.98	0.97
1:B:52:VAL:HG11	1:B:96:PHE:CE2	2.04	0.93
1:A:52:VAL:HG11	1:A:96:PHE:CZ	2.17	0.77
1:A:53:LYS:HE2	1:A:124:THR:HG21	1.83	0.60
1:B:52:VAL:HG11	1:B:96:PHE:CZ	2.38	0.59
1:A:53:LYS:HE2	1:A:124:THR:CG2	2.35	0.57
1:A:40:GLN:NE2	1:A:333:TYR:OH	2.43	0.51

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:56:PRO:HD3	2:B:402:NAG:C8	2.42	0.49
1:A:169:GLU:HG3	1:A:201:ARG:HA	1.94	0.48
1:B:169:GLU:HG3	1:B:201:ARG:HA	1.95	0.48
1:B:56:PRO:HD3	2:B:402:NAG:H81	1.94	0.48
1:B:236:ARG:HE	1:B:240:ALA:HB2	1.78	0.47
1:A:52:VAL:CG1	1:A:96:PHE:CZ	2.95	0.47
1:B:50:VAL:HG22	1:B:127:TRP:CE2	2.50	0.46
1:A:137:VAL:HB	1:A:213:PHE:HB3	1.98	0.45
1:B:47:LYS:HA	1:B:50:VAL:HG23	1.98	0.44
1:A:50:VAL:CG2	1:A:125:LEU:HB3	2.47	0.44
1:B:165:GLN:HG2	1:B:167:TYR:CZ	2.53	0.43
1:B:253:ARG:NH2	2:B:402:NAG:O6	2.51	0.43
1:B:121:ALA:HB3	1:B:256:LEU:HD12	2.01	0.41
1:A:50:VAL:HG21	1:A:125:LEU:HB3	2.01	0.41
1:B:147:THR:HG22	1:B:148:VAL:HG23	2.02	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	318/353 (90%)	313 (98%)	5 (2%)	0	100	100
1	B	311/353 (88%)	302 (97%)	9 (3%)	0	100	100
All	All	629/706 (89%)	615 (98%)	14 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	278/311 (89%)	275 (99%)	3 (1%)	65	79
1	B	275/311 (88%)	266 (97%)	9 (3%)	33	45
All	All	553/622 (89%)	541 (98%)	12 (2%)	45	61

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	160	THR
1	A	251	ASP
1	A	334	LEU
1	B	72	GLU
1	B	79	ASN
1	B	103	LEU
1	B	147	THR
1	B	153	LYS
1	B	186	SER
1	B	198	GLU
1	B	233	LEU
1	B	334	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	40	GLN
1	B	290	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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 2 are monoatomic - leaving 4 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	$\# Z > 2$
2	NAG	A	401	1	14,14,15	0.38	0	17,19,21	1.28	3 (17%)
2	NAG	A	402	1	14,14,15	0.55	0	17,19,21	1.03	1 (5%)
2	NAG	B	401	1	14,14,15	0.49	0	17,19,21	1.25	3 (17%)
2	NAG	B	402	1	14,14,15	0.48	0	17,19,21	1.01	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
2	NAG	A	401	1	-	0/6/23/26	0/1/1/1
2	NAG	A	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	B	402	1	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	401	NAG	C2-N2-C7	-2.57	119.45	122.90
2	A	401	NAG	C4-C3-C2	-2.54	107.30	111.02
2	B	401	NAG	C2-N2-C7	-2.43	119.64	122.90
2	B	401	NAG	C4-C3-C2	-2.33	107.60	111.02
2	B	401	NAG	C8-C7-N2	2.30	119.94	116.12

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	401	NAG	C8-C7-N2	2.24	119.83	116.12
2	A	402	NAG	C6-C5-C4	-2.14	107.76	113.02

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	402	NAG	3	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	322/353 (91%)	0.36	18 (5%) 30 27	8, 23, 47, 69	0
1	B	317/353 (89%)	0.32	20 (6%) 26 23	11, 23, 47, 60	0
All	All	639/706 (90%)	0.34	38 (5%) 28 25	8, 23, 47, 69	0

All (38) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	29	ASP	4.6
1	B	74	PRO	4.6
1	B	75	TYR	4.4
1	A	75	TYR	4.3
1	B	73	ASN	4.0
1	B	202	TRP	4.0
1	A	181	ARG	3.9
1	A	28	THR	3.8
1	A	202	TRP	3.7
1	B	183	ARG	3.6
1	A	27	THR	3.6
1	A	345	ALA	3.6
1	B	31	GLY	3.3
1	A	26	VAL	3.3
1	B	24	SER	3.2
1	B	345	ALA	3.1
1	B	263	GLY	3.0
1	B	25	TRP	2.9
1	A	183	ARG	2.9
1	B	262	GLY	2.8
1	B	265	TYR	2.8
1	A	178	SER	2.6
1	B	338	HIS	2.6
1	A	160	THR	2.6

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	B	264	THR	2.6
1	B	32	PRO	2.5
1	B	261	LEU	2.5
1	A	73	ASN	2.3
1	A	318	LYS	2.3
1	A	176	GLY	2.3
1	B	318	LYS	2.2
1	B	33	THR	2.2
1	B	72	GLU	2.2
1	A	25	TRP	2.2
1	A	124	THR	2.2
1	A	159	ARG	2.1
1	A	71	HIS	2.1
1	B	34	TRP	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($\text{\AA}^2$)	Q<0.9
2	NAG	A	402	14/15	0.80	0.15	36,40,46,48	0
2	NAG	B	402	14/15	0.81	0.14	33,38,41,43	0
2	NAG	A	401	14/15	0.94	0.08	12,15,18,20	0
2	NAG	B	401	14/15	0.97	0.06	11,14,17,17	0
3	CA	B	403	1/1	0.97	0.03	14,14,14,14	0
3	CA	A	403	1/1	0.99	0.01	15,15,15,15	0

6.5 Other polymers [i](#)

There are no such residues in this entry.