



wwPDB X-ray Structure Validation Summary Report ⓘ

Apr 17, 2017 – 01:04 PM EDT

PDB ID : 5B4X
Title : Crystal structure of the ApoER2 ectodomain in complex with the Reelin R56 fragment
Authors : Yasui, N.; Hirai, H.; Yamashita, K.; Takagi, J.; Nogi, T.
Deposited on : 2016-04-20
Resolution : 3.20 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20029077
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
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-20029077

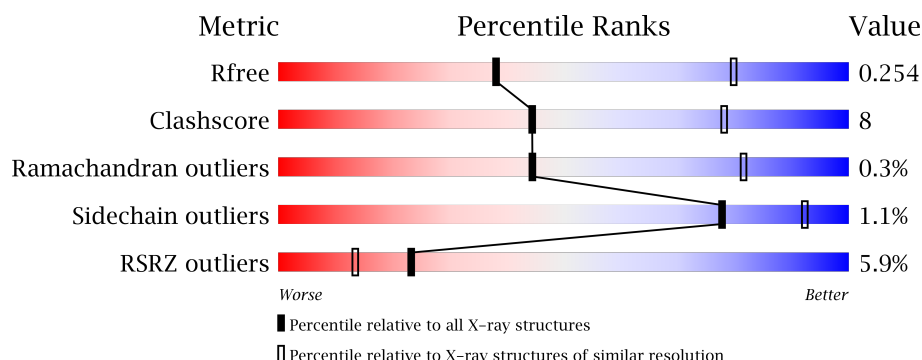
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.20 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}	100719	1015 (3.22-3.18)
Clashscore	112137	1009 (3.20-3.20)
Ramachandran outliers	110173	1118 (3.22-3.18)
Sidechain outliers	110143	1117 (3.22-3.18)
RSRZ outliers	101464	1020 (3.22-3.18)

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	725	<div> <div>0.1%</div> <div>82%</div> <div>16%</div> <div>0.1%</div> </div>
1	C	725	<div> <div>3%</div> <div>79%</div> <div>18%</div> <div>0.1%</div> </div>
2	B	570	<div> <div>9%</div> <div>67%</div> <div>16%</div> <div>0.1%</div> <div>16%</div> </div>
2	D	570	<div> <div>9%</div> <div>69%</div> <div>15%</div> <div>0.1%</div> <div>16%</div> </div>

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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	CA	C	4004	-	-	-	X
3	CA	D	5002	-	-	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 18892 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 Reelin.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	710	Total	C	N	O	S	Se	0	0	0
			5609	3559	955	1059	26	10			
1	C	709	Total	C	N	O	S	Se	0	0	0
			5601	3555	953	1057	26	10			

There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1946	GLY	-	expression tag	UNP Q60841
A	1947	ARG	-	expression tag	UNP Q60841
A	2101	ALA	CYS	engineered mutation	UNP Q60841
A	2663	ARG	-	expression tag	UNP Q60841
A	2664	LEU	-	expression tag	UNP Q60841
A	2665	GLU	-	expression tag	UNP Q60841
A	2666	ASN	-	expression tag	UNP Q60841
A	2667	LEU	-	expression tag	UNP Q60841
A	2668	TYR	-	expression tag	UNP Q60841
A	2669	PHE	-	expression tag	UNP Q60841
A	2670	GLN	-	expression tag	UNP Q60841
C	1946	GLY	-	expression tag	UNP Q60841
C	1947	ARG	-	expression tag	UNP Q60841
C	2101	ALA	CYS	engineered mutation	UNP Q60841
C	2663	ARG	-	expression tag	UNP Q60841
C	2664	LEU	-	expression tag	UNP Q60841
C	2665	GLU	-	expression tag	UNP Q60841
C	2666	ASN	-	expression tag	UNP Q60841
C	2667	LEU	-	expression tag	UNP Q60841
C	2668	TYR	-	expression tag	UNP Q60841
C	2669	PHE	-	expression tag	UNP Q60841
C	2670	GLN	-	expression tag	UNP Q60841

- Molecule 2 is a protein called Low density lipoprotein receptor-related protein 8, apolipoprotein e receptor, isoform CRA_e.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	479	Total 3798	C 2359	N 656	O 746	S 37	0	0	0
2	D	479	Total 3798	C 2359	N 656	O 746	S 37	0	0	0

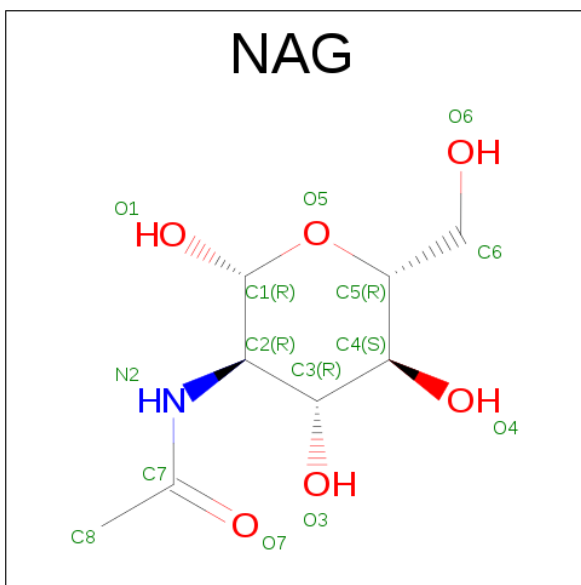
There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	40	SER	-	expression tag	UNP D3DQ39
B	607	ALA	-	expression tag	UNP D3DQ39
B	608	SER	-	expression tag	UNP D3DQ39
B	609	SER	-	expression tag	UNP D3DQ39
D	40	SER	-	expression tag	UNP D3DQ39
D	607	ALA	-	expression tag	UNP D3DQ39
D	608	SER	-	expression tag	UNP D3DQ39
D	609	SER	-	expression tag	UNP D3DQ39

- Molecule 3 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	4	Total 4	Ca 4	0	0
3	A	4	Total 4	Ca 4	0	0
3	D	4	Total 4	Ca 4	0	0
3	C	4	Total 4	Ca 4	0	0

- Molecule 4 is N-ACETYL-D-GLUCOSAMINE (three-letter code: NAG) (formula: C₈H₁₅NO₆).

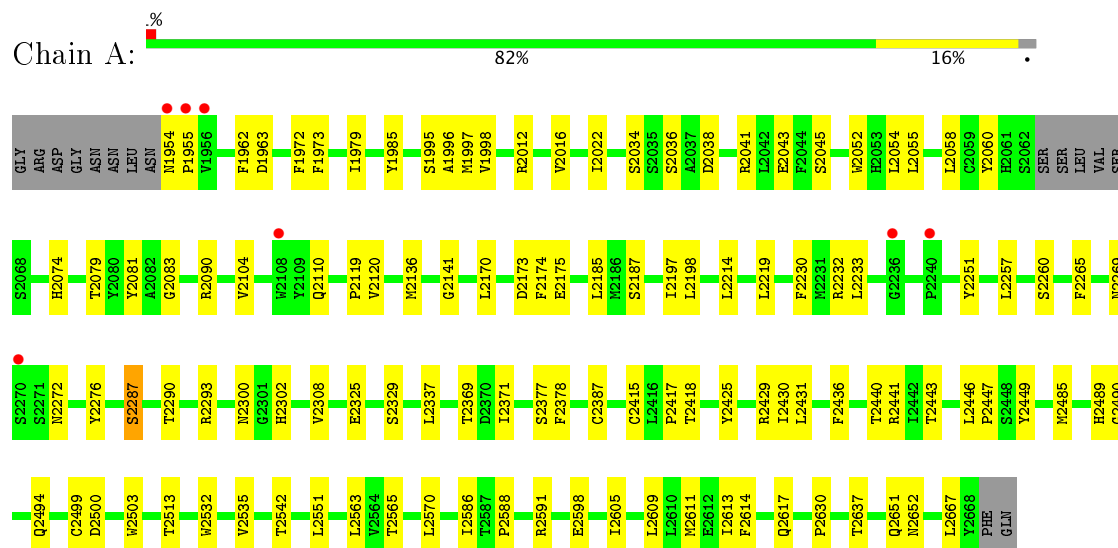


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			14	8	1	5		
4	A	1	Total	C	N	O	0	0
			14	8	1	5		
4	B	1	Total	C	N	O	0	0
			14	8	1	5		
4	C	1	Total	C	N	O	0	0
			14	8	1	5		
4	D	1	Total	C	N	O	0	0
			14	8	1	5		

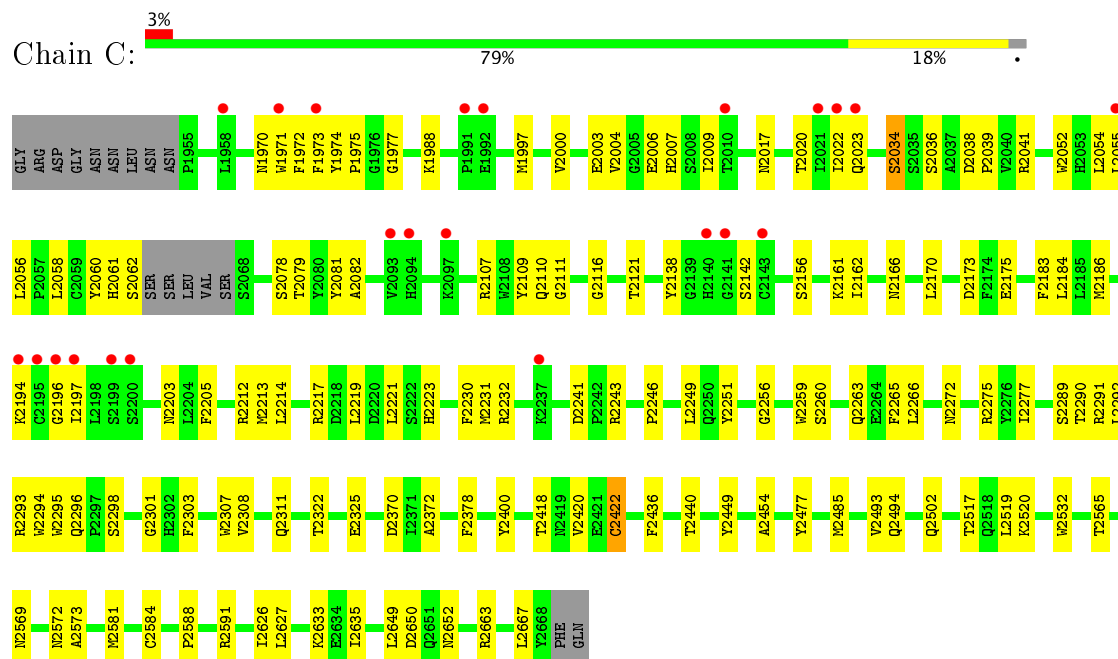
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA 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.

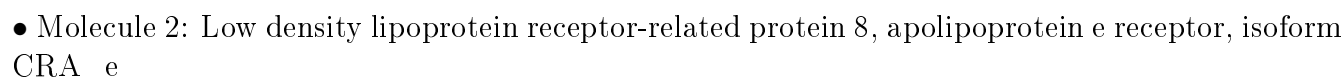
• Molecule 1: Reelin



• Molecule 1: Reelin



- Chain B:  9% 67% 16% 8%



4 Data and refinement statistics

Property	Value	Source
Space group	P 65	Depositor
Cell constants a, b, c, α , β , γ	205.95Å 205.95Å 169.84Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	49.52 – 3.20 49.47 – 3.20	Depositor EDS
% Data completeness (in resolution range)	100.0 (49.52-3.20) 100.0 (49.47-3.20)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.49 (at 3.19Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.200 , 0.255 0.199 , 0.254	Depositor DCC
R_{free} test set	3467 reflections (5.42%)	DCC
Wilson B-factor (Å ²)	82.6	Xtriage
Anisotropy	0.120	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 49.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.026 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	18892	wwPDB-VP
Average B, all atoms (Å ²)	99.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.74% 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, 0.333 respectively for untwinned datasets, and 0.375, 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: CA, 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	0.36	0/5755	0.57	0/7815
1	C	0.36	0/5747	0.58	0/7803
2	B	0.35	0/3885	0.60	0/5265
2	D	0.35	0/3885	0.59	0/5265
All	All	0.36	0/19272	0.58	0/26148

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
2	B	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	434	LYS	Peptide

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	5609	0	5322	68	0
1	C	5601	0	5318	94	0
2	B	3798	0	3572	67	0
2	D	3798	0	3572	57	0
3	A	4	0	0	0	0
3	B	4	0	0	0	0
3	C	4	0	0	0	0
3	D	4	0	0	0	0
4	A	28	0	26	0	0
4	B	14	0	13	0	0
4	C	14	0	13	1	0
4	D	14	0	13	0	0
All	All	18892	0	17849	277	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 8.

The worst 5 of 277 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:C:1975:PRO:HB2	1:C:2006:GLU:O	1.53	1.09
1:C:2023:GLN:NE2	1:C:2142:SER:HB2	1.79	0.97
2:D:367:HIS:ND1	2:D:387:GLY:HA3	1.81	0.95
1:C:2060:TYR:HE1	1:C:2062:SER:HG	1.13	0.92
1:C:2183:PHE:HB2	1:C:2186:MSE:HE2	1.56	0.87

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	706/725 (97%)	628 (89%)	76 (11%)	2 (0%)	44 81

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	705/725 (97%)	641 (91%)	64 (9%)	0	100	100
2	B	475/570 (83%)	431 (91%)	41 (9%)	3 (1%)	28	72
2	D	475/570 (83%)	421 (89%)	53 (11%)	1 (0%)	51	86
All	All	2361/2590 (91%)	2121 (90%)	234 (10%)	6 (0%)	44	81

5 of 6 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	2174	PHE
2	B	318	MET
2	B	550	GLU
2	B	592	PRO
2	D	592	PRO

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	621/624 (100%)	614 (99%)	7 (1%)	78	92
1	C	620/624 (99%)	617 (100%)	3 (0%)	91	97
2	B	427/497 (86%)	418 (98%)	9 (2%)	59	85
2	D	427/497 (86%)	422 (99%)	5 (1%)	75	91
All	All	2095/2242 (93%)	2071 (99%)	24 (1%)	78	92

5 of 24 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	B	318	MET
2	B	467	ARG
2	D	540	ASN
2	B	365	GLN
2	B	446	GLN

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

Mol	Chain	Res	Type
1	A	2572	ASN
1	C	2166	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 ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 21 ligands modelled in this entry, 16 are monoatomic - leaving 5 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	$\# Z > 2$
4	NAG	A	4005	1	14,14,15	0.47	0	15,19,21	1.49	2 (13%)
4	NAG	A	4006	1	14,14,15	0.56	0	15,19,21	1.31	2 (13%)
4	NAG	B	5005	2	14,14,15	0.46	0	15,19,21	1.91	1 (6%)
4	NAG	C	4005	1	14,14,15	0.51	0	15,19,21	1.54	2 (13%)
4	NAG	D	5005	2	14,14,15	0.58	0	15,19,21	1.70	2 (13%)

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
4	NAG	A	4005	1	-	0/6/23/26	0/1/1/1
4	NAG	A	4006	1	-	0/6/23/26	0/1/1/1
4	NAG	B	5005	2	-	0/6/23/26	0/1/1/1
4	NAG	C	4005	1	-	0/6/23/26	0/1/1/1
4	NAG	D	5005	2	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

The worst 5 of 9 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	5005	NAG	C6-C5-C4	-2.68	106.72	113.00
4	A	4005	NAG	O5-C1-C2	-2.20	108.41	111.47
4	A	4006	NAG	O5-C1-C2	-2.11	108.54	111.47
4	C	4005	NAG	O5-C1-C2	-2.07	108.60	111.47
4	A	4006	NAG	C1-O5-C5	3.36	116.80	112.17

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
4	C	4005	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	700/725 (96%)	0.05	7 (1%) 82 72	50, 89, 145, 176	0
1	C	699/725 (96%)	0.13	23 (3%) 47 31	44, 93, 158, 196	0
2	B	479/570 (84%)	0.42	54 (11%) 6 4	45, 79, 223, 273	0
2	D	479/570 (84%)	0.58	54 (11%) 6 4	51, 90, 232, 316	0
All	All	2357/2590 (91%)	0.26	138 (5%) 23 13	44, 88, 190, 316	0

The worst 5 of 138 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	111	PRO	15.7
2	D	112	ASP	12.6
2	D	94	ASN	11.4
2	B	122	THR	10.9
2	D	120	THR	9.4

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 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, 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(\AA^2)	Q<0.9
3	CA	C	4004	1/1	0.96	0.28	3.22	56,56,56,56	0
3	CA	A	4004	1/1	0.94	0.22	0.63	62,62,62,62	0
3	CA	C	4003	1/1	0.94	0.24	0.61	68,68,68,68	0
4	NAG	C	4005	14/15	0.86	0.21	0.55	77,82,93,101	0
4	NAG	A	4006	14/15	0.89	0.28	0.52	80,93,104,108	0
4	NAG	A	4005	14/15	0.93	0.21	0.20	105,112,121,123	0
4	NAG	B	5005	14/15	0.88	0.23	0.16	74,83,92,93	0
3	CA	A	4003	1/1	0.93	0.20	-0.04	66,66,66,66	0
3	CA	A	4002	1/1	0.99	0.21	-0.10	118,118,118,118	0
3	CA	A	4001	1/1	0.86	0.13	-0.84	108,108,108,108	0
3	CA	D	5004	1/1	0.90	0.07	-1.42	129,129,129,129	0
3	CA	C	4001	1/1	0.82	0.10	-1.49	120,120,120,120	0
3	CA	D	5002	1/1	0.20	0.49	-1.64	306,306,306,306	0
3	CA	B	5003	1/1	0.63	0.08	-1.64	196,196,196,196	0
3	CA	C	4002	1/1	0.81	0.07	-1.82	132,132,132,132	0
3	CA	D	5003	1/1	0.83	0.09	-2.14	213,213,213,213	0
3	CA	B	5002	1/1	0.72	0.38	-2.22	276,276,276,276	0
3	CA	B	5004	1/1	0.94	0.06	-2.35	113,113,113,113	0
3	CA	B	5001	1/1	0.98	0.06	-2.89	86,86,86,86	0
3	CA	D	5001	1/1	0.91	0.05	-4.07	97,97,97,97	0
4	NAG	D	5005	14/15	0.88	0.20	-	79,85,89,101	0

6.5 Other polymers [i](#)

There are no such residues in this entry.