



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 5, 2024 – 04:00 pm BST

PDB ID : 2WJW
Title : Crystal structure of the human ionotropic glutamate receptor GluR2 ATD region at 1.8 Å resolution
Authors : Clayton, A.; Siebold, C.; Gilbert, R.J.C.; Sutton, G.C.; Harlos, K.; McIlhinney, R.A.J.; Jones, E.Y.; Aricescu, A.R.
Deposited on : 2009-06-01
Resolution : 1.80 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

<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	:	4.02b-467
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	3.0
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.003 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.39

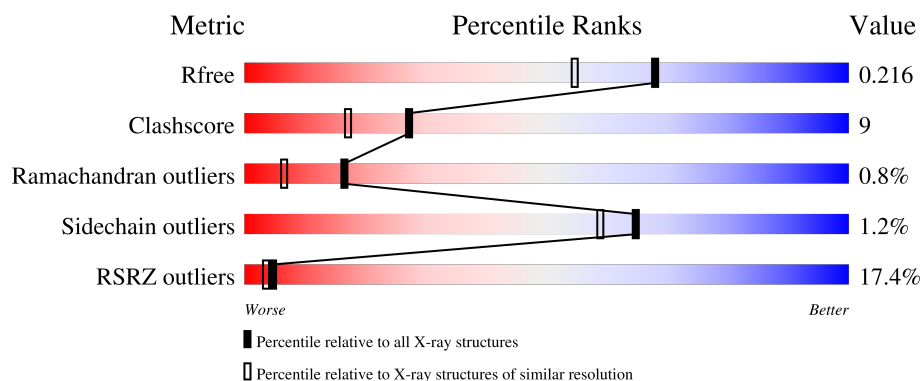
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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}	164625	7108 (1.80-1.80)
Clashscore	180529	8162 (1.80-1.80)
Ramachandran outliers	177936	8077 (1.80-1.80)
Sidechain outliers	177891	8076 (1.80-1.80)
RSRZ outliers	164620	7108 (1.80-1.80)

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 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	388	<div> <div>16%</div> <div>79%</div> <div>15%</div> <div>• 5%</div> </div>

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 3294 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 GLUTAMATE RECEPTOR 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	368	Total	C	N	O	S	0	7	0
			2991	1904	507	570	10			

- Molecule 2 is 2-acetamido-2-deoxy-beta-D-glucopyranose (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		

- Molecule 3 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

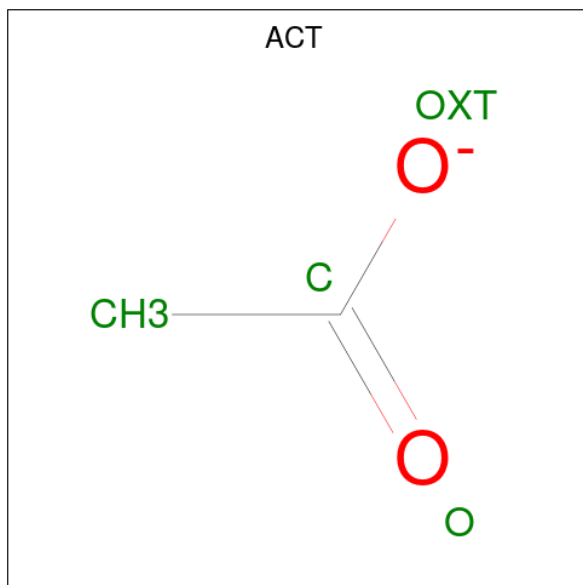
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Cl	0	0
			1	1		

- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	S	0	0
			5	4	1		

- Molecule 5 is ACETATE ION (three-letter code: ACT) (formula: $\text{C}_2\text{H}_3\text{O}_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			4	2	2		

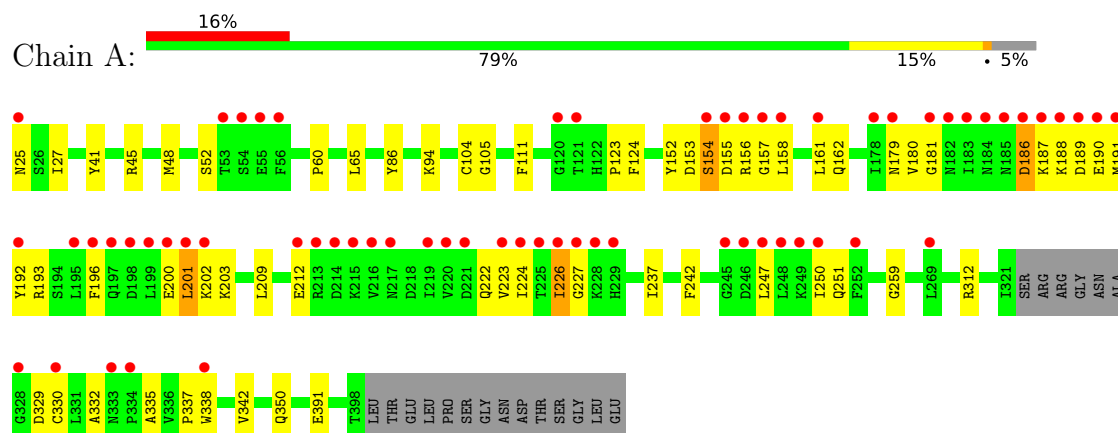
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	279	Total 279	O 279	0	0

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: GLUTAMATE RECEPTOR 2



4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	50.98Å 122.45Å 138.85Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.16 – 1.80 39.16 – 1.80	Depositor EDS
% Data completeness (in resolution range)	95.7 (39.16-1.80) 99.9 (39.16-1.80)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.93 (at 1.79Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, R_{free}	0.181 , 0.217 0.183 , 0.216	Depositor DCC
R_{free} test set	1637 reflections (4.02%)	wwPDB-VP
Wilson B-factor (Å ²)	20.6	Xtriage
Anisotropy	0.678	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 51.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	3294	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.14% 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: NAG, CL, SO4, ACT

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	0.30	0/3052	0.68	2/4126 (0.0%)

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	1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	186	ASP	CB-CG-OD2	-28.39	92.75	118.30
1	A	186	ASP	OD1-CG-OD2	-14.11	96.49	123.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	186	ASP	Sidechain

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	2991	0	2938	54	0
2	A	14	0	13	0	0
3	A	1	0	0	0	0
4	A	5	0	0	0	0
5	A	4	0	3	1	0
6	A	279	0	0	7	0
All	All	3294	0	2954	54	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 9.

All (54) 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:391:GLU:HG3	6:A:2252:HOH:O	1.75	0.87
1:A:154[A]:SER:HB3	1:A:179:ASN:HD21	1.43	0.80
1:A:154[A]:SER:HB3	1:A:179:ASN:ND2	2.03	0.72
1:A:27:ILE:HG23	1:A:86:TYR:HD2	1.56	0.70
1:A:162:GLN:HG3	6:A:2106:HOH:O	1.89	0.70
1:A:94:LYS:HD3	1:A:155[B]:ASP:HA	1.74	0.68
1:A:154[A]:SER:H	1:A:181:GLY:HA3	1.61	0.64
1:A:201:LEU:HD12	1:A:201:LEU:C	2.23	0.59
1:A:180:VAL:HG12	1:A:180:VAL:O	2.02	0.58
1:A:350:GLN:HG3	6:A:2228:HOH:O	2.04	0.56
1:A:188:LYS:HD3	1:A:189:ASP:N	2.21	0.55
1:A:329:ASP:HB3	1:A:332:ALA:HB2	1.88	0.54
1:A:187:LYS:HD2	1:A:187:LYS:H	1.73	0.54
1:A:247:LEU:HD22	6:A:2141:HOH:O	2.07	0.54
1:A:48:MET:SD	1:A:60:PRO:HG3	2.47	0.53
1:A:201:LEU:C	1:A:203:LYS:H	2.12	0.52
1:A:187:LYS:O	1:A:191:MET:HG3	2.11	0.50
1:A:192:TYR:HB3	1:A:222:GLN:HG2	1.94	0.49
1:A:196:PHE:O	1:A:200:GLU:HB2	2.13	0.49
1:A:391:GLU:HB3	6:A:2273:HOH:O	2.13	0.49
1:A:154[A]:SER:OG	1:A:181:GLY:HA3	2.14	0.48
1:A:86:TYR:HD1	1:A:338:TRP:CH2	2.32	0.48
1:A:41:TYR:CE2	1:A:45:ARG:HD2	2.49	0.47
1:A:222:GLN:O	1:A:226:ILE:HG22	2.15	0.47
1:A:312:ARG:NH1	6:A:2206:HOH:O	2.48	0.47
1:A:48:MET:O	1:A:52:SER:HB3	2.14	0.47
1:A:188:LYS:HG2	1:A:192:TYR:CE2	2.50	0.47
1:A:187:LYS:HD2	1:A:187:LYS:N	2.30	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:196:PHE:CE1	1:A:223:VAL:HG22	2.51	0.46
1:A:156[A]:ARG:HG2	1:A:157[A]:GLY:H	1.80	0.46
1:A:65:LEU:HD23	1:A:65:LEU:N	2.30	0.45
1:A:202:LYS:HG2	1:A:202:LYS:O	2.17	0.45
1:A:226:ILE:O	1:A:226:ILE:HG23	2.17	0.45
1:A:158[B]:LEU:HD12	1:A:158[B]:LEU:HA	1.72	0.44
1:A:153[B]:ASP:OD2	1:A:212:GLU:HG3	2.17	0.44
1:A:190:GLU:OE1	1:A:193:ARG:HG3	2.17	0.44
1:A:335:ALA:O	1:A:337:PRO:HD3	2.18	0.43
1:A:154[A]:SER:HB3	1:A:179:ASN:CG	2.39	0.43
1:A:250:ILE:HG22	1:A:251:GLN:N	2.34	0.43
1:A:153[B]:ASP:O	1:A:154[B]:SER:HB2	2.18	0.43
1:A:342:VAL:HG23	6:A:2223:HOH:O	2.18	0.43
1:A:104:CYS:SG	1:A:111:PHE:HB2	2.58	0.43
1:A:242:PHE:CD1	1:A:259:GLY:HA3	2.54	0.43
1:A:152:TYR:CE1	1:A:158[B]:LEU:HD11	2.54	0.42
1:A:158[B]:LEU:HG	1:A:161:LEU:HD23	2.00	0.42
1:A:158[B]:LEU:O	1:A:161:LEU:HB3	2.19	0.42
1:A:27:ILE:HG23	1:A:86:TYR:CD2	2.46	0.42
1:A:154[A]:SER:N	1:A:181:GLY:HA3	2.33	0.41
1:A:105:GLY:C	5:A:1402:ACT:H2	2.40	0.41
1:A:123:PRO:O	1:A:124:PHE:HB2	2.20	0.41
1:A:242:PHE:O	1:A:247:LEU:HD23	2.21	0.41
1:A:209:LEU:HD12	1:A:237:ILE:CD1	2.51	0.41
1:A:153[B]:ASP:O	1:A:154[B]:SER:CB	2.69	0.40
1:A:27:ILE:HD12	1:A:27:ILE:N	2.37	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	371/388 (96%)	352 (95%)	15 (4%)	4 (1%)	12 3

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	227	GLY
1	A	154[A]	SER
1	A	154[B]	SER
1	A	226	ILE

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	326/336 (97%)	322 (99%)	4 (1%)	67 62

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	25	ASN
1	A	201	LEU
1	A	224	ILE
1	A	330	CYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	25	ASN

5.3.3 RNA ⓘ

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](#)

Of 4 ligands modelled in this entry, 1 is monoatomic - leaving 3 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	SO4	A	1401	-	4,4,4	0.11	0	6,6,6	0.13	0
2	NAG	A	1399	1	14,14,15	0.50	0	17,19,21	1.03	1 (5%)
5	ACT	A	1402	-	3,3,3	0.76	0	3,3,3	1.34	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	1399	1	-	1/6/23/26	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	1399	NAG	C1-O5-C5	2.53	115.62	112.19

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1399	NAG	C4-C5-C6-O6

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	1402	ACT	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, 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	368/388 (94%)	0.44	64 (17%) 5 3	11, 25, 104, 357	8 (2%)

All (64) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	157[A]	GLY	8.5
1	A	201	LEU	7.7
1	A	183	ILE	6.0
1	A	247	LEU	5.2
1	A	155[A]	ASP	4.4
1	A	154[A]	SER	4.4
1	A	186	ASP	4.1
1	A	226	ILE	4.0
1	A	192	TYR	3.9
1	A	191	MET	3.9
1	A	184	ASN	3.9
1	A	219	ILE	3.8
1	A	187	LYS	3.8
1	A	245	GLY	3.7
1	A	328	GLY	3.6
1	A	330	CYS	3.6
1	A	213	ARG	3.5
1	A	182	ASN	3.4
1	A	249	LYS	3.4
1	A	199	LEU	3.3
1	A	224	ILE	3.3
1	A	188	LYS	3.2
1	A	250	ILE	3.2
1	A	246	ASP	3.2
1	A	56	PHE	3.2
1	A	185	ASN	3.1
1	A	338	TRP	3.1

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	158[A]	LEU	3.1
1	A	161	LEU	3.0
1	A	216	VAL	3.0
1	A	223	VAL	3.0
1	A	200	GLU	2.9
1	A	248	LEU	2.9
1	A	195	LEU	2.8
1	A	156[A]	ARG	2.8
1	A	220	VAL	2.8
1	A	212	GLU	2.8
1	A	227	GLY	2.6
1	A	198	ASP	2.6
1	A	202	LYS	2.6
1	A	120	GLY	2.6
1	A	214	ASP	2.6
1	A	53	THR	2.5
1	A	178	ILE	2.5
1	A	228	LYS	2.4
1	A	225	THR	2.4
1	A	215	LYS	2.4
1	A	121	THR	2.4
1	A	229	HIS	2.4
1	A	181	GLY	2.4
1	A	54	SER	2.3
1	A	217	ASN	2.3
1	A	252	PHE	2.3
1	A	334	PRO	2.2
1	A	25	ASN	2.2
1	A	221	ASP	2.2
1	A	179	ASN	2.2
1	A	333	ASN	2.2
1	A	190	GLU	2.2
1	A	55	GLU	2.1
1	A	197	GLN	2.1
1	A	189	ASP	2.0
1	A	269	LEU	2.0
1	A	196	PHE	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 monosaccharides 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(\AA^2)	Q<0.9
2	NAG	A	1399	14/15	0.80	0.13	26,39,54,55	0
5	ACT	A	1402	4/4	0.81	0.18	55,60,62,63	0
3	CL	A	1400	1/1	0.85	0.11	90,90,90,90	0
4	SO4	A	1401	5/5	0.95	0.10	32,34,41,50	0

6.5 Other polymers [i](#)

There are no such residues in this entry.