



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 06:17 AM GMT

PDB ID : 2WJW
Title : CRYSTAL STRUCTURE OF THE HUMAN IONOTROPIC GLUTAMATE RECEPTOR GLUR2 ATD REGION AT 1.8 Å RESOLUTION
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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
<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 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
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) : trunk26865

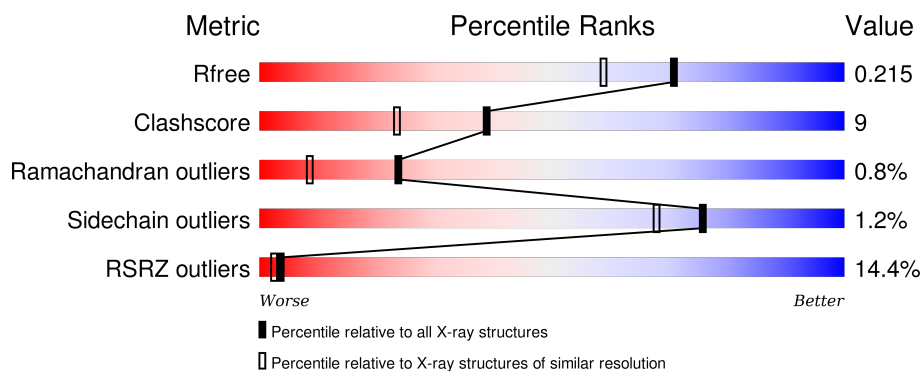
1 Overall quality at a glance

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}	91344	4533 (1.80-1.80)
Clashscore	102246	5383 (1.80-1.80)
Ramachandran outliers	100387	5320 (1.80-1.80)
Sidechain outliers	100360	5319 (1.80-1.80)
RSRZ outliers	91569	4547 (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 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	388	

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 criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	NAG	A	1399	-	-	-	X

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 SUGAR (N-ACETYL-D-GLUCOSAMINE) (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).

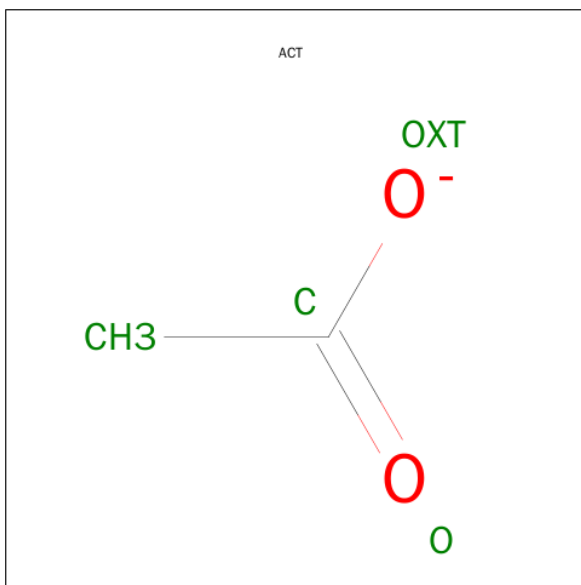
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Cl	0	0
			1	1		

- 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: $\text{C}_2\text{H}_3\text{O}_2$).

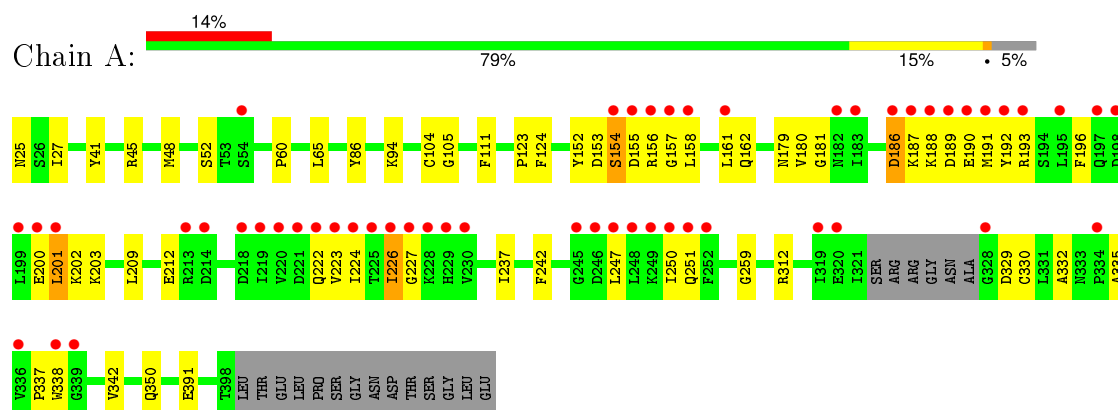


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			4	2	2		

- Molecule 6 is water.

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

• 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 47.06 – 1.80	Depositor EDS
% Data completeness (in resolution range)	95.7 (39.16-1.80) 99.7 (47.06-1.80)	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 , 0.215	Depositor DCC
R_{free} test set	1637 reflections (4.02%)	DCC
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 , 57.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 40754 reflections	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, 0.375 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: CL, NAG, 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

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:162:GLN:HG3	6:A:2106:HOH:O	1.89	0.70
1:A:27:ILE:HG23	1:A:86:TYR:HD2	1.56	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:C	1:A:201:LEU:HD12	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:H	1:A:187:LYS:HD2	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

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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:O	1:A:202:LYS:HG2	2.17	0.45
1:A:226:ILE:HG23	1:A:226:ILE:O	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:242:PHE:CD1	1:A:259:GLY:HA3	2.54	0.43
1:A:104:CYS:SG	1:A:111:PHE:HB2	2.58	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:N	1:A:27:ILE:HD12	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%)	17 5

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%)	78 71

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

Some 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 carbohydrates 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	$\# Z > 2$
2	NAG	A	1399	1	14,14,15	0.48	0	15,19,21	1.10	1 (6%)
4	SO4	A	1401	-	4,4,4	0.19	0	6,6,6	0.08	0
5	ACT	A	1402	-	1,3,3	1.23	0	0,3,3	0.00	-

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	-	0/6/23/26	0/1/1/1
4	SO4	A	1401	-	-	0/0/0/0	0/0/0/0
5	ACT	A	1402	-	-	0/0/0/0	0/0/0/0

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	2.66	115.62	112.25

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
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%)	0.52	53 (14%) 3 2	11, 26, 104, 378	1 (0%)

All (53) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	224	ILE	7.4
1	A	201	LEU	7.3
1	A	157[A]	GLY	7.1
1	A	225	THR	7.1
1	A	226	ILE	6.3
1	A	247	LEU	6.0
1	A	228	LYS	5.6
1	A	246	ASP	5.6
1	A	249	LYS	4.9
1	A	336	VAL	4.8
1	A	193	ARG	4.5
1	A	220	VAL	4.5
1	A	190	GLU	4.4
1	A	192	TYR	4.3
1	A	200	GLU	4.3
1	A	223	VAL	4.3
1	A	189	ASP	4.3
1	A	195	LEU	4.3
1	A	197	GLN	4.2
1	A	338	TRP	4.2
1	A	334	PRO	4.1
1	A	230	VAL	4.1
1	A	251	GLN	4.0
1	A	186	ASP	3.9
1	A	155[A]	ASP	3.9
1	A	219	ILE	3.8
1	A	156[A]	ARG	3.8

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Mol	Chain	Res	Type	RSRZ
1	A	221	ASP	3.7
1	A	250	ILE	3.7
1	A	227	GLY	3.6
1	A	187	LYS	3.5
1	A	229	HIS	3.4
1	A	183	ILE	3.4
1	A	182	ASN	3.4
1	A	158[A]	LEU	3.3
1	A	154[A]	SER	3.3
1	A	191	MET	3.2
1	A	54	SER	3.2
1	A	214	ASP	3.1
1	A	198	ASP	3.1
1	A	245	GLY	2.9
1	A	213	ARG	2.7
1	A	161	LEU	2.6
1	A	199	LEU	2.5
1	A	222	GLN	2.3
1	A	188	LYS	2.3
1	A	252	PHE	2.3
1	A	320	GLU	2.2
1	A	328	GLY	2.1
1	A	339	GLY	2.1
1	A	319	ILE	2.0
1	A	218	ASP	2.0
1	A	248	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 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
2	NAG	A	1399	14/15	0.72	0.23	3.59	26,39,54,55	0
4	SO4	A	1401	5/5	0.98	0.12	0.12	32,34,41,50	0
5	ACT	A	1402	4/4	0.83	0.17	-	55,60,62,63	0
3	CL	A	1400	1/1	0.92	0.09	-	90,90,90,90	0

6.5 Other polymers [i](#)

There are no such residues in this entry.