



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 13, 2020 – 11:49 AM BST

PDB ID : 3OEN
Title : Crystal structure of GluN2D ligand-binding core in complex with L-glutamate
Authors : Simorowski, N.; Furukawa, H.
Deposited on : 2010-08-12
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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Xtriage (Phenix) : 1.13
EDS : 2.14.4.dev1
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.14.4.dev1

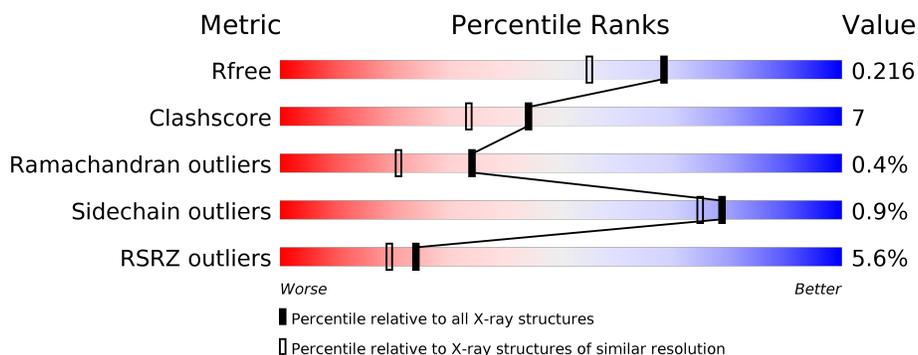
1 Overall quality at a glance

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}	130704	5950 (1.80-1.80)
Clashscore	141614	6793 (1.80-1.80)
Ramachandran outliers	138981	6697 (1.80-1.80)
Sidechain outliers	138945	6696 (1.80-1.80)
RSRZ outliers	127900	5850 (1.80-1.80)

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 2277 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.

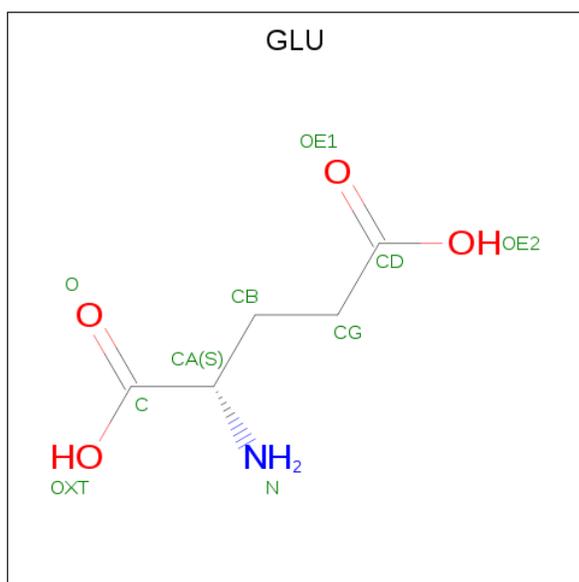
- Molecule 1 is a protein called Glutamate [NMDA] receptor subunit epsilon-4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	250	1954	1243	337	360	14	0	0	0

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	GLY	-	EXPRESSION TAG	UNP Q62645
A	143	GLY	-	LINKER	UNP Q62645
A	144	THR	-	LINKER	UNP Q62645

- Molecule 2 is GLUTAMIC ACID (three-letter code: GLU) (formula: C₅H₉NO₄).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
2	A	1	10	5	1	4	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	313	Total 313	O 313	0	0

SEQUENCE-PLOTS INFOmissingINFO

3 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	60.07Å 113.70Å 95.33Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.42 – 1.80 28.65 – 1.80	Depositor EDS
% Data completeness (in resolution range)	98.0 (19.42-1.80) 99.7 (28.65-1.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.49 (at 1.80Å)	Xtrriage
Refinement program	PHENIX (phenix.refine)	Depositor
R, R_{free}	0.186 , 0.217 0.185 , 0.216	Depositor DCC
R_{free} test set	1538 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	24.1	Xtrriage
Anisotropy	0.276	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 58.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	2277	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.88% 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.

4 Model quality [i](#)

4.1 Standard geometry [i](#)

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	0.35	0/1992	0.55	0/2687

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

4.2 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	1954	0	1949	26	0
2	A	10	0	5	1	0
3	A	313	0	0	9	0
All	All	2277	0	1954	26	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 7.

All (26) 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:38:ARG:HH12	1:A:79:ASP:CG	1.94	0.71
1:A:267:LEU:HG	1:A:272:ILE:HD12	1.73	0.70
1:A:36:PRO:HB3	1:A:55:GLU:H	1.54	0.70

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:199:GLU:HB2	3:A:298:HOH:O	1.98	0.64
1:A:38:ARG:NH1	1:A:79:ASP:OD1	2.33	0.61
1:A:238:LYS:HD2	1:A:239:VAL:N	2.16	0.61
1:A:56:LYS:HB3	3:A:580:HOH:O	2.04	0.58
1:A:199:GLU:HG2	3:A:584:HOH:O	2.03	0.58
1:A:163:PRO:HB2	1:A:188:TYR:OH	2.04	0.58
1:A:203:GLN:OE1	3:A:311:HOH:O	2.17	0.57
1:A:199:GLU:CG	3:A:584:HOH:O	2.53	0.56
1:A:116:THR:HG1	2:A:1001:GLU:N	2.04	0.55
1:A:36:PRO:HB3	1:A:55:GLU:N	2.22	0.54
1:A:4:THR:N	3:A:301:HOH:O	2.42	0.53
1:A:196:ARG:HG3	3:A:298:HOH:O	2.08	0.53
1:A:55:GLU:O	1:A:56:LYS:HB3	2.12	0.49
1:A:164:LEU:O	1:A:188:TYR:HE2	1.96	0.47
1:A:214:TYR:HB3	1:A:219:LEU:HG	1.97	0.47
1:A:257:ARG:HB3	1:A:258:PRO:HD3	1.96	0.46
1:A:124:ILE:HD12	1:A:125:VAL:HG13	1.98	0.46
1:A:56:LYS:HD2	1:A:56:LYS:C	2.36	0.45
1:A:277:ARG:HD3	3:A:565:HOH:O	2.17	0.45
1:A:178:ILE:HG22	1:A:186:HIS:HB2	1.99	0.45
1:A:164:LEU:O	1:A:188:TYR:CE2	2.71	0.44
1:A:32:ARG:NH1	3:A:387:HOH:O	2.49	0.43
1:A:233:THR:OG1	1:A:237:GLY:HA2	2.22	0.40

There are no symmetry-related clashes.

4.3 Torsion angles [\(i\)](#)

4.3.1 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	244/286 (85%)	231 (95%)	12 (5%)	1 (0%)	34 21

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	56	LYS

4.3.2 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	211/246 (86%)	209 (99%)	2 (1%)	78 75

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	238	LYS
1	A	251	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	106	GLN
1	A	193	ASN
1	A	194	GLN
1	A	250	HIS

4.3.3 RNA [i](#)

There are no RNA molecules in this entry.

4.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

4.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

4.6 Ligand geometry

1 ligand is modelled in this entry.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

4.7 Other polymers

There are no such residues in this entry.

4.8 Polymer linkage issues

There are no chain breaks in this entry.

5 Fit of model and data [i](#)

5.1 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	250/286 (87%)	0.26	14 (5%) 24 19	15, 26, 46, 72	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	54	PRO	9.7
1	A	236	SER	6.8
1	A	4	THR	6.6
1	A	55	GLU	6.2
1	A	239	VAL	3.9
1	A	163	PRO	3.8
1	A	240	PHE	3.8
1	A	235	GLY	3.3
1	A	238	LYS	2.9
1	A	285	HIS	2.8
1	A	237	GLY	2.7
1	A	234	ILE	2.6
1	A	17	PHE	2.4
1	A	114	SER	2.0

5.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.4 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	GLU	A	1001	10/10	0.97	0.14	14,16,20,21	0

5.5 Other polymers [i](#)

There are no such residues in this entry.