

Full wwPDB X-ray Structure Validation Report (i)

May 16, 2020 – 02:13 pm BST

PDB ID : 5IKB

Title: Crystal structure of the kainate receptor GluK4 ligand binding domain in

complex with kainate

Authors: Kristensen, O.; Kristensen, L.B.; Frydenvang, K.; Kastrup, J.S.

Deposited on : 2016-03-03

Resolution : 2.05 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 (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

MolProbity : 4.02b-467

Mogul: 1.8.5 (274361), CSD as541be (2020)

Xtriage (Phenix) : 1.13 EDS : 2.11

Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)

Refmac: 5.8.0158

CCP4 : 7.0.044 (Gargrove) roteins) : Engh & Huber (2001)

Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)

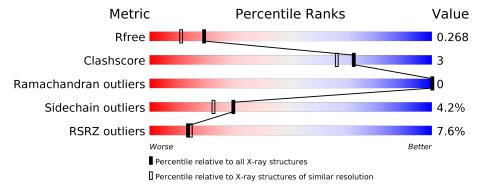
Validation Pipeline (wwPDB-VP) : 2.11

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: $X\text{-}RAY\ DIFFRACTION$

The reported resolution of this entry is 2.05 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	$\begin{array}{c} \text{Whole archive} \\ (\#\text{Entries}) \end{array}$	$\begin{array}{c} {\rm Similar \; resolution} \\ (\#{\rm Entries, \; resolution \; range(\AA)}) \end{array}$
R_{free}	130704	1692 (2.04-2.04)
Clashscore	141614	1773 (2.04-2.04)
Ramachandran outliers	138981	1752 (2.04-2.04)
Sidechain outliers	138945	1752 (2.04-2.04)
RSRZ outliers	127900	1672 (2.04-2.04)

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 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 <=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		
			7%		
1	A	257	85%	11%	•



2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 2292 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.

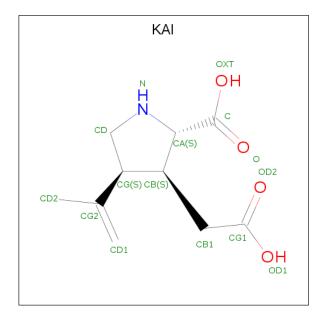
• Molecule 1 is a protein called Glutamate receptor ionotropic, kainate 4, Glutamate receptor ionotropic, kainate 4.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	A	249	Total	С	N	О	S	0	7	0
		_ 10	2045	1298	342	393	12		•	

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	GLY	-	expression tag	UNP Q01812
A	2	PRO	1	expression tag	UNP Q01812
A	115	ARG	-	$_{ m linker}$	UNP Q01812
A	116	LYS	-	linker	UNP Q01812
A	117	GLY	-	$_{ m linker}$	UNP Q01812
A	118	THR	-	$_{ m linker}$	UNP Q01812

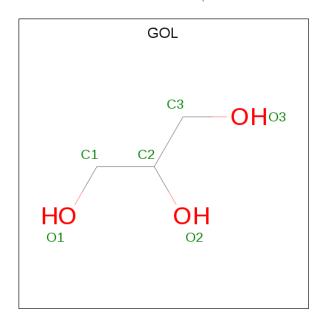
• Molecule 2 is 3-(CARBOXYMETHYL)-4-ISOPROPENYLPROLINE (three-letter code: KAI) (formula: C₁₀H₁₅NO₄).





Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total 15	C 10	N 1	O 4	0	0

• Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 6 3 3	0	0
3	A	1	Total C O 6 3 3	0	0

• Molecule 4 is water.

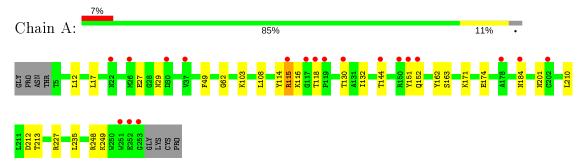
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	220	Total O 220 220	0	0



3 Residue-property plots (i)

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: Glutamate receptor ionotropic, kainate 4, Glutamate receptor ionotropic, kainate 4





4 Data and refinement statistics (i)

Property	Value	Source	
Space group	C 2 2 21	Depositor	
Cell constants	44.50Å 69.87Å 162.76Å	Depositor	
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor	
Resolution (Å)	40.69 - 2.05	Depositor	
Resolution (A)	40.69 - 2.05	EDS	
% Data completeness	99.1 (40.69-2.05)	Depositor	
(in resolution range)	99.1 (40.69-2.05)	EDS	
R_{merge}	0.14	Depositor	
R_{sym}	(Not available)	Depositor	
$< I/\sigma(I) > 1$	1.67 (at 2.05Å)	Xtriage	
Refinement program	REFMAC 5.8.0135, PHENIX 1.10.1	Depositor	
D D	0.191 , 0.262	Depositor	
R, R_{free}	0.199 , 0.268	DCC	
R_{free} test set	897 reflections (5.53%)	wwPDB-VP	
Wilson B-factor (Å ²)	23.9	Xtriage	
Anisotropy	0.119	Xtriage	
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.36 , 54.9	EDS	
L-test for twinning ²	$< L >=0.46, < L^2>=0.29$	Xtriage	
Estimated twinning fraction	No twinning to report.	Xtriage	
F_o, F_c correlation	0.93	EDS	
Total number of atoms	2292	wwPDB-VP	
Average B, all atoms (\mathring{A}^2)	27.0	wwPDB-VP	

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 5.77% of the height of the origin peak. No significant pseudotranslation is detected.

²Theoretical values of <|L|>, $< L^2>$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.



¹Intensities estimated from amplitudes.

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: GOL, KAI

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	Boı	nd lengths	Bo	ond angles
IVIOI	Chain	RMSZ	# Z > 5	RMSZ	# Z > 5
1	A	0.89	1/2084~(0.0%)	0.92	3/2816 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	${ m Observed}({ m \AA})$	$\operatorname{Ideal}(ext{\AA})$
1	A	62	GLY	C-O	5.03	1.31	1.23

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^o)$	$Ideal(^{o})$
1	A	248	ARG	NE-CZ-NH1	6.67	123.64	120.30
1	A	212	ASP	CB-CG-OD1	5.28	123.05	118.30
1	A	227	ARG	NE-CZ-NH1	5.12	122.86	120.30

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	2045	0	2005	14	1
2	A	15	0	13	0	0
3	A	12	0	16	2	0
4	A	220	0	0	6	0

Continued on next page...



Continued from previous page...

Mol	Chain	Non-H	H(model)	$\mathbf{H}(\mathbf{added})$	Clashes	Symm-Clashes
All	All	2292	0	2034	14	1

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 3.

All (14) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic	Clash
Atom-1	Atom-2	${f distance}\;({ m \AA})$	${ m overlap} ({ m \AA})$
1:A:103:LYS:HE3	1:A:235:LEU:HD11	1.86	0.57
1:A:174[A]:GLU:OE1	1:A:174[A]:GLU:HA	2.07	0.55
1:A:118:THR:HG22	4:A:486:HOH:O	2.06	0.54
1:A:49:PHE:O	3:A:303:GOL:C1	2.59	0.51
1:A:171:LYS:NZ	4:A:407:HOH:O	2.44	0.49
1:A:29:ASN:HB3	4:A:596:HOH:O	2.14	0.47
1:A:115:ARG:NH2	1:A:184:ASN:OD1	2.49	0.46
1:A:213[A]:THR:HG23	1:A:213[A]:THR:O	2.15	0.46
1:A:249:LYS:NZ	4:A:409:HOH:O	2.45	0.46
1:A:116:LYS:HA	4:A:486:HOH:O	2.16	0.46
1:A:174[A]:GLU:HG2	4:A:541:HOH:O	2.18	0.44
1:A:12:LEU:HD23	1:A:17:LEU:HD12	2.00	0.43
1:A:144:THR:HG21	3:A:302:GOL:H31	2.00	0.43
1:A:114:TYR:OH	1:A:132:ILE:HD12	2.19	0.42

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	$egin{array}{c} ext{Interatomic} \ ext{distance} \ (ext{Å}) \end{array}$	$egin{array}{c} ext{Clash} \ ext{overlap } (ext{Å}) \end{array}$
1:A:162:TYR:OH	1:A:184:ASN:ND2[8 456]	2.06	0.14

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	$254/257 \ (99\%)$	245 (96%)	9 (4%)	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	221/220 (100%)	211 (96%)	10 (4%)	27 20

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	27	GLU
1	A	108	LEU
1	A	115	ARG
1	A	130	THR
1	A	151	TYR
1	A	152	GLN
1	A	163[A]	SER
1	A	163[B]	SER
1	A	201	ASN
1	A	210	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA (i)

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)

3 ligands are modelled in this entry.

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		Res	Tinle	Link Bond lengths			В	ond ang	gles	
MIGI	Type	Chain	nes	LIIIK	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	GOL	A	302	_	5,5,5	0.29	0	5,5,5	0.72	0
3	GOL	A	303	_	5,5,5	0.49	0	5,5,5	0.78	0
2	KAI	A	301	-	6,15,15	0.86	0	8,21,21	0.88	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	${f Res}$	Link	Chirals	${f Torsions}$	Rings
3	GOL	A	302	_	-	2/4/4/4	-
3	GOL	A	303	-	-	3/4/4/4	-
2	KAI	A	301	-	-	4/6/25/25	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	303	GOL	O1-C1-C2-C3
2	A	301	KAI	CD-CG-CG2-CD1
3	A	303	GOL	C1-C2-C3-O3
3	A	303	GOL	O1-C1-C2-O2

Continued on next page...



 $Continued\ from\ previous\ page...$

Mol	Chain	Res	Type	Atoms
2	A	301	KAI	CD-CG-CG2-CD2
2	A	301	KAI	CB-CG-CG2-CD2
2	A	301	KAI	CB-CG-CG2-CD1
3	A	302	GOL	O2-C2-C3-O3
3	A	302	GOL	O1-C1-C2-C3

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	302	GOL	1	0
3	A	303	GOL	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, 95^{th} 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 $>$	$\#\mathrm{RSRZ}{>}2$	$OWAB(A^2)$	Q < 0.9
1	A	249/257 (96%)	0.62	19 (7%) 13 14	14, 23, 49, 73	0

All (19) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	151	TYR	4.9
1	A	202	CYS	4.1
1	A	130	THR	3.9
1	A	30	ASP	3.1
1	A	26	MET	3.0
1	A	184	ASN	2.8
1	A	119	PRO	2.8
1	A	178	ALA	2.7
1	A	150	ARG	2.7
1	A	22	ASN	2.6
1	A	117	GLY	2.5
1	A	118	THR	2.4
1	A	252	GLU	2.4
1	A	115	ARG	2.3
1	A	251	TRP	2.3
1	A	253	GLY	2.3
1	A	37	VAL	2.3
1	A	144	THR	2.2
1	A	152	GLN	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. The B-factors column lists the minimum, median, 95^{th} 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	${f Res}$	Atoms	RSCC	RSR	${f B-factors({ m \AA}^2)}$	Q<0.9
3	GOL	A	303	6/6	0.74	0.16	43,50,52,56	0
3	GOL	A	302	6/6	0.89	0.15	25,27,28,30	0
2	KAI	A	301	15/15	0.92	0.14	15,18,23,24	0

6.5 Other polymers (i)

There are no such residues in this entry.

