



wwPDB X-ray Structure Validation Summary Report ⓘ

Oct 15, 2017 – 06:05 PM EDT

PDB ID : 3BDB
Title : Crystal Structure of Novel Immune-Type Receptor 11 Extracellular Fragment from *Ictalurus punctatus* including Stalk Region
Authors : Ostrov, D.A.; Hernandez Prada, J.A.; Haire, R.N.; Cannon, J.P.; Magis, A.T.; Bailey, K.M.; Litman, G.W.
Deposited on : unknown
Resolution : 2.80 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
Xtriage (Phenix) : 1.9-1692
EDS : rb-20030345
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-20030345

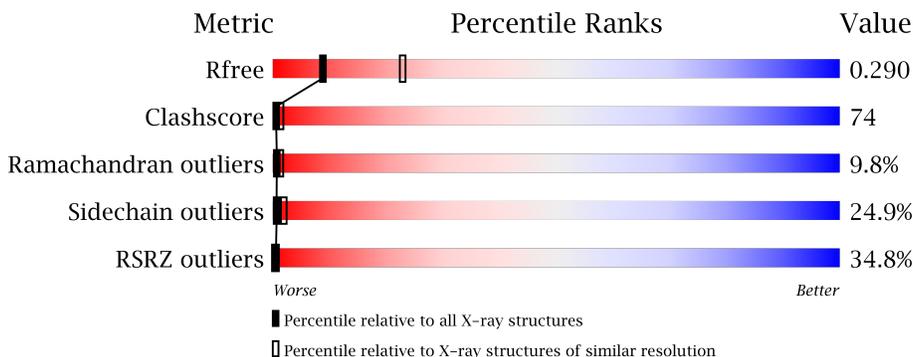
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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}	100719	2583 (2.80-2.80)
Clashscore	112137	3033 (2.80-2.80)
Ramachandran outliers	110173	2983 (2.80-2.80)
Sidechain outliers	110143	2985 (2.80-2.80)
RSRZ outliers	101464	2610 (2.80-2.80)

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	137	
1	B	137	
1	C	137	
1	D	137	
1	E	137	

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Mol	Chain	Length	Quality of chain
1	F	137	 <p>A horizontal bar chart representing the quality of chain. The bar is divided into five segments with the following percentages from left to right: 16% (green), 28% (red), 42% (yellow), 29% (orange), and 8% (grey). A small black dot is located on the orange segment.</p>

2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 6021 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 Novel immune-type receptor 11.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	126	995	620	170	199	6	0	0	0
1	B	126	995	620	170	199	6	0	0	0
1	C	126	995	620	170	199	6	0	0	0
1	D	126	995	620	170	199	6	0	0	0
1	E	126	995	620	170	199	6	0	0	0
1	F	126	995	620	170	199	6	0	0	0

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	INITIATING METHIONINE	UNP Q8UWK4
B	1	MET	-	INITIATING METHIONINE	UNP Q8UWK4
C	1	MET	-	INITIATING METHIONINE	UNP Q8UWK4
D	1	MET	-	INITIATING METHIONINE	UNP Q8UWK4
E	1	MET	-	INITIATING METHIONINE	UNP Q8UWK4
F	1	MET	-	INITIATING METHIONINE	UNP Q8UWK4

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	9	Total	O	0	0
			9	9		
2	B	11	Total	O	0	0
			11	11		
2	C	8	Total	O	0	0
			8	8		

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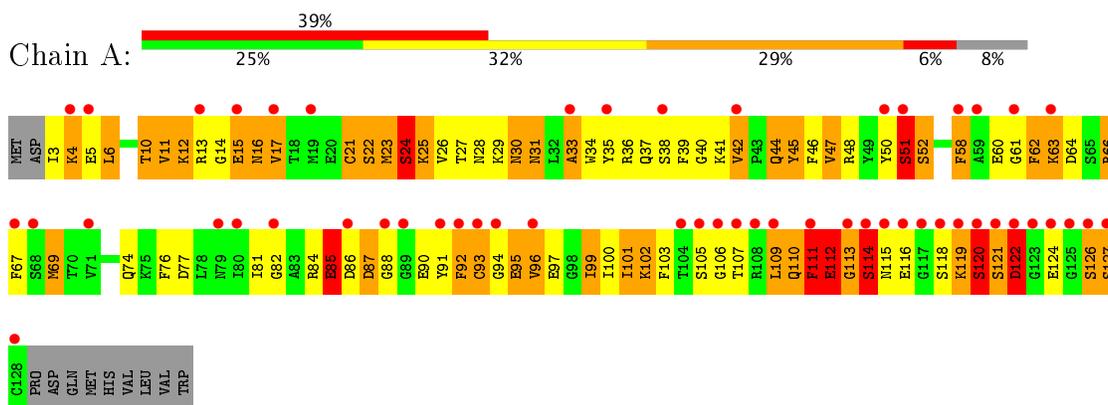
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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	D	9	Total O 9 9	0	0
2	E	7	Total O 7 7	0	0
2	F	7	Total O 7 7	0	0

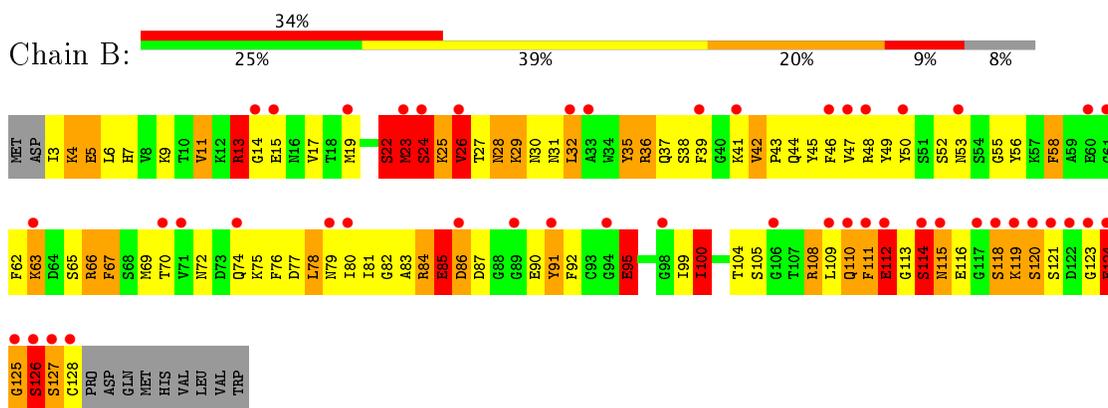
3 Residue-property plots [i](#)

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.

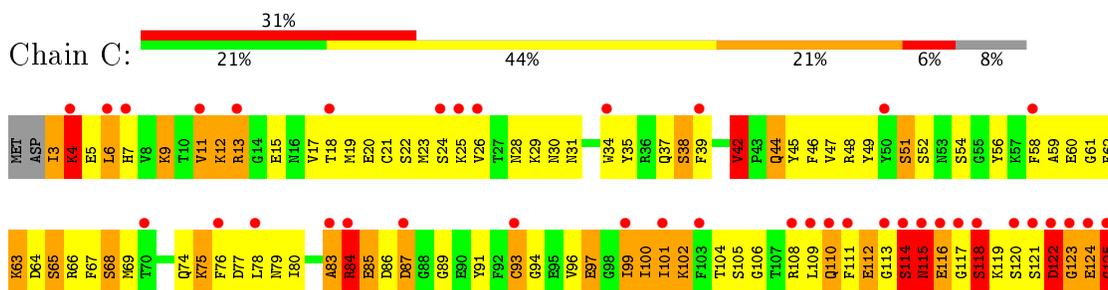
- Molecule 1: Novel immune-type receptor 11



- Molecule 1: Novel immune-type receptor 11

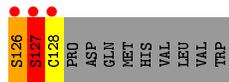
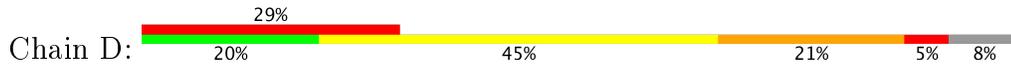


- Molecule 1: Novel immune-type receptor 11

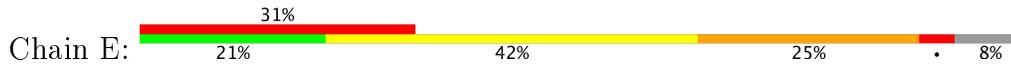




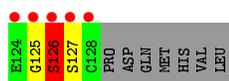
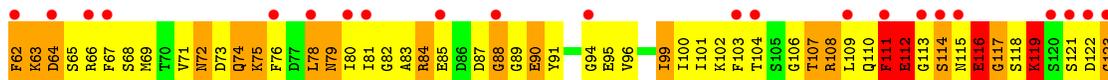
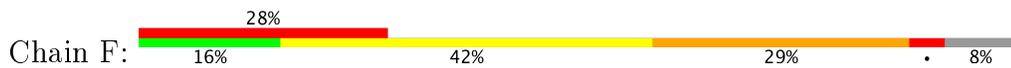
● Molecule 1: Novel immune-type receptor 11



● Molecule 1: Novel immune-type receptor 11



● Molecule 1: Novel immune-type receptor 11



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	91.58Å 91.34Å 115.33Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.80 29.51 – 2.81	Depositor EDS
% Data completeness (in resolution range)	98.1 (30.00-2.80) 99.3 (29.51-2.81)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	27.26 (at 2.80Å)	Xtrriage
Refinement program	CNS	Depositor
R, R_{free}	0.284 , 0.291 0.286 , 0.290	Depositor DCC
R_{free} test set	1231 reflections (5.14%)	DCC
Wilson B-factor (Å ²)	66.1	Xtrriage
Anisotropy	0.258	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 39.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	0.427 for k,h,-l	Xtrriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	6021	wwPDB-VP
Average B, all atoms (Å ²)	70.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.06% 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](#)

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	1.30	8/1013 (0.8%)	1.13	7/1353 (0.5%)
1	B	1.35	7/1013 (0.7%)	1.12	4/1353 (0.3%)
1	C	1.43	12/1013 (1.2%)	1.20	4/1353 (0.3%)
1	D	1.46	10/1013 (1.0%)	1.25	9/1353 (0.7%)
1	E	1.46	5/1013 (0.5%)	1.22	4/1353 (0.3%)
1	F	1.65	11/1013 (1.1%)	1.40	7/1353 (0.5%)
All	All	1.45	53/6078 (0.9%)	1.22	35/8118 (0.4%)

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	13
1	B	0	13
1	C	0	13
1	D	0	13
1	E	0	14
1	F	0	8
All	All	0	74

The worst 5 of 53 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	90	GLU	CD-OE2	5.99	1.32	1.25
1	A	76	PHE	CD2-CE2	5.99	1.51	1.39
1	F	35	TYR	CD1-CE1	5.98	1.48	1.39
1	D	103	PHE	CE1-CZ	5.90	1.48	1.37
1	F	71	VAL	CA-CB	-5.90	1.42	1.54

The worst 5 of 35 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	83	ALA	N-CA-CB	-5.99	101.72	110.10
1	F	36	ARG	NE-CZ-NH1	5.89	123.25	120.30
1	C	9	LYS	N-CA-C	-5.84	95.22	111.00
1	E	23	MET	CG-SD-CE	-5.83	90.87	100.20
1	D	18	THR	OG1-CB-CG2	-5.79	96.69	110.00

There are no chirality outliers.

5 of 74 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	10	THR	Peptide
1	A	22	SER	Peptide
1	A	23	MET	Peptide
1	A	33	ALA	Peptide
1	A	82	GLY	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	995	0	945	151	0
1	B	995	0	945	166	0
1	C	995	0	945	181	0
1	D	995	0	947	130	0
1	E	995	0	945	166	0
1	F	995	0	945	149	0
2	A	9	0	0	2	0
2	B	11	0	0	7	0
2	C	8	0	0	1	0
2	D	9	0	0	4	0
2	E	7	0	0	2	0
2	F	7	0	0	7	0
All	All	6021	0	5672	858	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 74.

The worst 5 of 858 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:F:62:PHE:CZ	1:F:64:ASP:HB2	1.66	1.30
1:C:12:LYS:HA	1:C:112:GLU:CG	1.61	1.29
1:B:126:SER:HB2	1:B:127:SER:CA	1.65	1.21
1:C:85:GLU:HA	1:C:111:PHE:CZ	1.75	1.20
1:A:99:ILE:HG21	1:B:31:ASN:ND2	1.54	1.19

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	124/137 (90%)	91 (73%)	19 (15%)	14 (11%)	0	1
1	B	124/137 (90%)	96 (77%)	18 (14%)	10 (8%)	1	2
1	C	124/137 (90%)	87 (70%)	23 (18%)	14 (11%)	0	1
1	D	124/137 (90%)	92 (74%)	24 (19%)	8 (6%)	1	4
1	E	124/137 (90%)	85 (68%)	27 (22%)	12 (10%)	1	1
1	F	124/137 (90%)	88 (71%)	21 (17%)	15 (12%)	0	1
All	All	744/822 (90%)	539 (72%)	132 (18%)	73 (10%)	1	1

5 of 73 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	15	GLU
1	A	51	SER
1	A	85	GLU
1	A	115	ASN
1	A	121	SER

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	109/120 (91%)	83 (76%)	26 (24%)	1	2
1	B	109/120 (91%)	81 (74%)	28 (26%)	0	1
1	C	109/120 (91%)	86 (79%)	23 (21%)	1	3
1	D	109/120 (91%)	80 (73%)	29 (27%)	0	1
1	E	109/120 (91%)	79 (72%)	30 (28%)	0	1
1	F	109/120 (91%)	82 (75%)	27 (25%)	1	2
All	All	654/720 (91%)	491 (75%)	163 (25%)	1	2

5 of 163 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	114	SER
1	D	78	LEU
1	F	65	SER
1	C	118	SER
1	D	23	MET

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 25 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	31	ASN
1	C	110	GLN
1	F	74	GLN
1	C	44	GLN
1	D	31	ASN

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](#)

There are no ligands in this entry.

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	126/137 (91%)	2.46	53 (42%) 0 0	32, 63, 100, 100	0
1	B	126/137 (91%)	2.11	47 (37%) 0 0	31, 62, 100, 100	0
1	C	126/137 (91%)	2.08	42 (33%) 0 0	33, 60, 100, 100	0
1	D	126/137 (91%)	2.11	40 (31%) 0 0	26, 61, 100, 100	0
1	E	126/137 (91%)	1.82	43 (34%) 0 0	48, 79, 100, 100	0
1	F	126/137 (91%)	1.78	38 (30%) 1 0	43, 80, 100, 100	0
All	All	756/822 (91%)	2.06	263 (34%) 0 0	26, 71, 100, 100	0

The worst 5 of 263 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	128	CYS	16.8
1	D	128	CYS	14.8
1	D	120	SER	13.5
1	C	121	SER	12.3
1	B	123	GLY	12.1

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](#)

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.