



wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 14, 2017 – 06:12 am GMT

PDB ID : 1JFM
Title : CRYSTAL STRUCTURE OF MURINE NK CELL LIGAND RAE-1 BETA
Authors : Li, P.; Strong, R.K.
Deposited on : 2001-06-21
Resolution : 2.85 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) : NOT EXECUTED
EDS : NOT EXECUTED
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : recalc28949

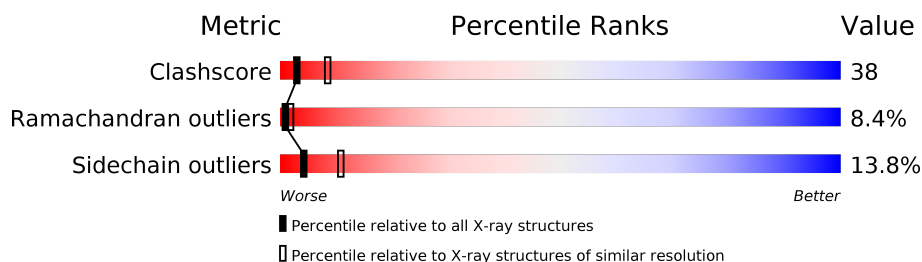
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.85 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))
Clashscore	112137	2749 (2.90-2.82)
Ramachandran outliers	110173	2687 (2.90-2.82)
Sidechain outliers	110143	2690 (2.90-2.82)

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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	174	
1	B	174	
1	C	174	
1	D	174	
1	E	174	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 7067 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.

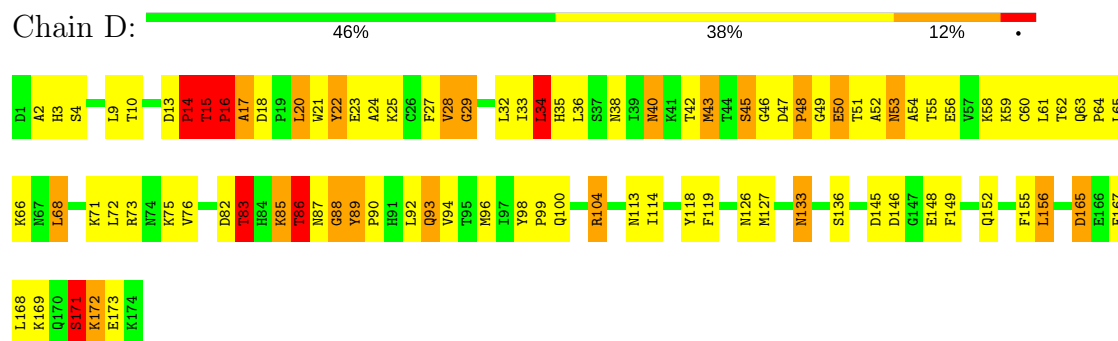
- Molecule 1 is a protein called retinoic acid early transcript beta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	174	Total	C	N	O	S	0	0	0
			1405	887	238	270	10			
1	B	174	Total	C	N	O	S	0	0	0
			1405	887	238	270	10			
1	C	174	Total	C	N	O	S	0	0	0
			1405	887	238	270	10			
1	D	174	Total	C	N	O	S	0	0	0
			1405	887	238	270	10			
1	E	174	Total	C	N	O	S	0	0	0
			1405	887	238	270	10			

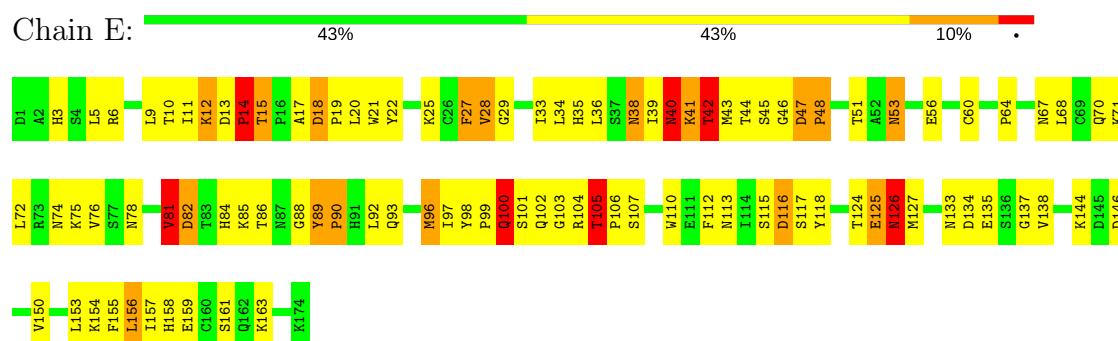
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	8	Total	O	0	0
			8	8		
2	B	6	Total	O	0	0
			6	6		
2	C	10	Total	O	0	0
			10	10		
2	D	12	Total	O	0	0
			12	12		
2	E	6	Total	O	0	0
			6	6		

● Molecule 1: retinoic acid early transcript beta



● Molecule 1: retinoic acid early transcript beta



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	I 41 2 2	Depositor
Cell constants a, b, c, α , β , γ	138.41Å 138.41Å 337.55Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	24.96 – 2.85	Depositor
% Data completeness (in resolution range)	96.9 (24.96-2.85)	Depositor
R_{merge}	0.08	Depositor
R_{sym}	0.08	Depositor
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.234 , 0.279	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	7067	wwPDB-VP
Average B, all atoms (Å ²)	89.0	wwPDB-VP

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	0.51	0/1438	0.83	3/1942 (0.2%)
1	B	0.49	0/1438	0.78	2/1942 (0.1%)
1	C	0.47	0/1438	0.83	5/1942 (0.3%)
1	D	0.59	0/1438	1.08	11/1942 (0.6%)
1	E	0.46	0/1438	0.83	4/1942 (0.2%)
All	All	0.51	0/7190	0.88	25/9710 (0.3%)

There are no bond length outliers.

The worst 5 of 25 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	D	40	ASN	N-CA-C	12.54	144.85	111.00
1	D	171	SER	N-CA-C	12.51	144.78	111.00
1	A	102	GLN	N-CA-C	-9.71	84.78	111.00
1	D	85	LYS	N-CA-C	8.74	134.61	111.00
1	D	86	THR	N-CA-C	8.20	133.14	111.00

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	1405	0	1375	106	1
1	B	1405	0	1375	102	0
1	C	1405	0	1375	117	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	1405	0	1375	95	0
1	E	1405	0	1375	115	0
2	A	8	0	0	1	0
2	B	6	0	0	0	0
2	C	10	0	0	1	0
2	D	12	0	0	0	0
2	E	6	0	0	0	0
All	All	7067	0	6875	526	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 38.

The worst 5 of 526 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:B:42:THR:O	1:B:43:MET:HG2	1.52	1.09
1:A:39:ILE:HG22	1:A:40:ASN:H	1.17	1.08
1:D:15:THR:HG22	1:D:16:PRO:HD2	1.23	1.08
1:C:15:THR:O	1:C:17:ALA:N	1.90	1.05
1:C:14:PRO:HB3	1:C:19:PRO:HA	1.40	0.99

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	Interatomic distance (Å)	Clash overlap (Å)
1:A:173:GLU:OE1	1:A:173:GLU:OE1[10_665]	2.19	0.01

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	172/174 (99%)	142 (83%)	16 (9%)	14 (8%)	1	2
1	B	172/174 (99%)	141 (82%)	23 (13%)	8 (5%)	3	8
1	C	172/174 (99%)	133 (77%)	25 (14%)	14 (8%)	1	2
1	D	172/174 (99%)	137 (80%)	17 (10%)	18 (10%)	0	1
1	E	172/174 (99%)	127 (74%)	27 (16%)	18 (10%)	0	1
All	All	860/870 (99%)	680 (79%)	108 (13%)	72 (8%)	1	2

5 of 72 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	2	ALA
1	A	15	THR
1	A	39	ILE
1	A	42	THR
1	A	47	ASP

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	160/160 (100%)	133 (83%)	27 (17%)	2	6
1	B	160/160 (100%)	146 (91%)	14 (9%)	12	31
1	C	160/160 (100%)	134 (84%)	26 (16%)	3	7
1	D	160/160 (100%)	141 (88%)	19 (12%)	6	16
1	E	160/160 (100%)	136 (85%)	24 (15%)	3	9
All	All	800/800 (100%)	690 (86%)	110 (14%)	4	11

5 of 110 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	47	ASP
1	C	134	ASP
1	E	84	HIS

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Mol	Chain	Res	Type
1	C	56	GLU
1	C	82	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 38 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	8	ASN
1	C	126	ASN
1	E	70	GLN
1	C	63	GLN
1	C	152	GLN

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

6.1 Protein, DNA and RNA chains

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.