



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 2, 2016 – 12:01 AM GMT

PDB ID : 8I1B
Title : A COMPARISON OF THE HIGH RESOLUTION STRUCTURES OF HUMAN AND MURINE INTERLEUKIN-1B
Authors : Ohlendorf, D.H.; Weber, P.C.; Salemme, F.R.
Deposited on : 1991-01-29
Resolution : 2.40 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) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

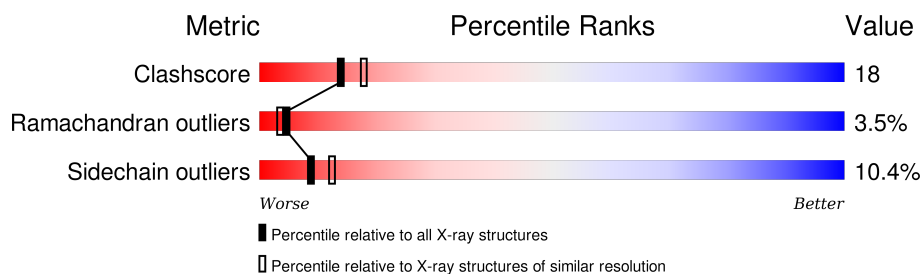
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	102246	3407 (2.40-2.40)
Ramachandran outliers	100387	3351 (2.40-2.40)
Sidechain outliers	100360	3352 (2.40-2.40)

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	152	 55% 28% 11% • •

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 1276 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 INTERLEUKIN-1 BETA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	146	Total	C	N	O	S	0	4	0
			1190	761	199	225	5			

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	86	Total	O	0	0
			86	86		

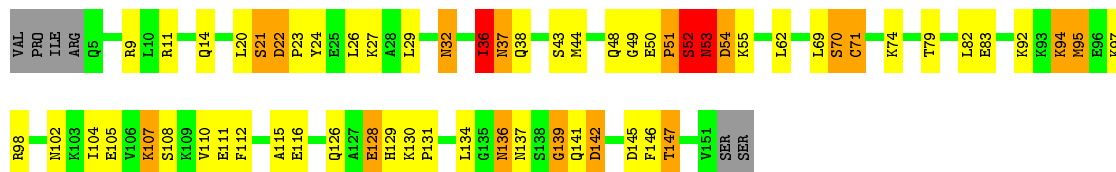
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

Note EDS was not executed.

• Molecule 1: INTERLEUKIN-1 BETA

Chain A: 



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 43	Depositor
Cell constants a, b, c, α , β , γ	55.71Å 55.71Å 79.00Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	5.00 – 2.40	Depositor
% Data completeness (in resolution range)	(Not available) (5.00-2.40)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	PROLSQ	Depositor
R, R_{free}	0.159 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	1276	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

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	1.18	1/1237 (0.1%)	1.83	24/1666 (1.4%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	70	SER	CA-CB	5.22	1.60	1.52

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	11	ARG	NE-CZ-NH2	-11.41	114.60	120.30
1	A	9	ARG	NE-CZ-NH2	-10.16	115.22	120.30
1	A	9	ARG	NE-CZ-NH1	9.63	125.12	120.30
1	A	147	THR	CA-CB-CG2	8.25	123.95	112.40
1	A	95	MET	CA-CB-CG	7.90	126.72	113.30
1	A	11	ARG	NE-CZ-NH1	7.88	124.24	120.30
1	A	70	SER	N-CA-CB	-7.78	98.83	110.50
1	A	43	SER	N-CA-CB	7.29	121.44	110.50
1	A	92	LYS	CA-CB-CG	6.77	128.30	113.40
1	A	27	LYS	N-CA-CB	6.66	122.59	110.60
1	A	36[A]	ILE	CB-CA-C	6.59	124.77	111.60
1	A	36[B]	ILE	CB-CA-C	6.59	124.77	111.60
1	A	54	ASP	CB-CG-OD1	6.33	123.99	118.30
1	A	128	GLU	OE1-CD-OE2	6.11	130.63	123.30
1	A	115	ALA	N-CA-CB	5.72	118.11	110.10
1	A	54	ASP	O-C-N	5.52	131.54	122.70
1	A	71	CYS	O-C-N	5.50	131.50	122.70
1	A	105	GLU	OE1-CD-OE2	5.39	129.77	123.30
1	A	32	ASN	CB-CA-C	5.36	121.12	110.40
1	A	142	ASP	CB-CG-OD1	-5.34	113.50	118.30
1	A	53	ASN	CA-C-N	-5.33	105.47	117.20
1	A	26	LEU	CA-CB-CG	5.28	127.44	115.30
1	A	11	ARG	N-CA-CB	5.25	120.05	110.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	21	SER	C-N-CA	5.03	134.28	121.70

There are no chirality outliers.

There are no planarity outliers.

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	1190	0	1195	44	0
2	A	86	0	0	9	0
All	All	1276	0	1195	44	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 18.

All (44) 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:50:GLU:HB2	1:A:51:PRO:CD	1.83	1.07
1:A:50:GLU:CB	1:A:51:PRO:HD2	1.90	1.01
1:A:50:GLU:HB2	1:A:51:PRO:HD2	0.96	0.93
1:A:137:ASN:ND2	2:A:365:HOH:O	1.98	0.90
1:A:107:LYS:O	1:A:108:SER:HB2	1.77	0.82
1:A:97:LYS:HG3	1:A:97:LYS:O	1.87	0.74
1:A:69:LEU:HD23	1:A:82:LEU:HD23	1.71	0.72
1:A:50:GLU:N	1:A:50:GLU:OE1	2.25	0.70
1:A:23:PRO:O	1:A:24:TYR:HB2	1.95	0.66
1:A:37:ASN:H	1:A:37:ASN:ND2	1.92	0.66
1:A:44:MET:HB2	2:A:400:HOH:O	1.96	0.65
1:A:97:LYS:HG2	1:A:116:GLU:OE1	2.00	0.61
1:A:111:GLU:OE2	1:A:145:ASP:OD1	2.19	0.61
1:A:14[A]:GLN:NE2	1:A:126:GLN:OE1	2.36	0.59
1:A:83:GLU:OE2	2:A:235:HOH:O	2.17	0.57
1:A:126:GLN:NE2	1:A:141:GLN:O	2.34	0.57
1:A:98:ARG:HG2	2:A:249:HOH:O	2.04	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:128:GLU:HG2	1:A:129:HIS:CD2	2.40	0.56
1:A:136:ASN:ND2	2:A:303:HOH:O	2.39	0.55
1:A:98:ARG:HG2	1:A:98:ARG:HH11	1.73	0.53
1:A:21:SER:C	1:A:23:PRO:HD2	2.30	0.52
1:A:110[A]:VAL:HG12	1:A:146:PHE:O	2.11	0.51
1:A:32:ASN:O	1:A:36[B]:ILE:HD13	2.11	0.50
1:A:107:LYS:O	1:A:108:SER:CB	2.48	0.50
1:A:130:LYS:HB3	1:A:131:PRO:HD2	1.93	0.50
1:A:126:GLN:HG3	1:A:141:GLN:HB3	1.93	0.49
1:A:49:GLY:C	1:A:50:GLU:OE1	2.51	0.49
1:A:49:GLY:HA2	2:A:325:HOH:O	2.14	0.48
1:A:131:PRO:HG2	2:A:232:HOH:O	2.14	0.47
1:A:52:SER:O	1:A:53:ASN:CB	2.63	0.47
1:A:71:CYS:HA	1:A:79:THR:O	2.15	0.47
1:A:55:LYS:HB3	1:A:102:ASN:HB3	1.99	0.45
1:A:69:LEU:HD23	1:A:69:LEU:HA	1.75	0.44
1:A:139:GLY:C	1:A:142:ASP:H	2.21	0.44
1:A:20:LEU:HD22	1:A:62:LEU:HD21	2.00	0.43
1:A:136:ASN:O	1:A:137:ASN:HB3	2.18	0.42
1:A:48:GLN:CD	1:A:94:LYS:HD2	2.40	0.42
1:A:137:ASN:N	2:A:216:HOH:O	2.48	0.42
1:A:74:LYS:HA	2:A:311:HOH:O	2.20	0.41
1:A:52:SER:O	1:A:53:ASN:HB2	2.20	0.41
1:A:110[B]:VAL:CG1	1:A:112:PHE:CE2	3.03	0.41
1:A:50:GLU:CB	1:A:51:PRO:CD	2.66	0.41
1:A:55:LYS:HG2	1:A:104:ILE:HG12	2.03	0.41

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	148/152 (97%)	134 (90%)	9 (6%)	5 (3%)	5 4

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	22	ASP
1	A	51	PRO
1	A	52	SER
1	A	53	ASN
1	A	139	GLY

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	139/141 (99%)	124 (89%)	15 (11%)	8 11

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	ASP
1	A	29	LEU
1	A	36[A]	ILE
1	A	36[B]	ILE
1	A	37	ASN
1	A	38	GLN
1	A	52	SER
1	A	54	ASP
1	A	70	SER
1	A	94	LYS
1	A	95	MET
1	A	107	LYS
1	A	134	LEU
1	A	136	ASN
1	A	147	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	34	GLN
1	A	35	ASN
1	A	37	ASN
1	A	38	GLN
1	A	129	HIS
1	A	136	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 ⓘ

6.1 Protein, DNA and RNA chains ⓘ

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates ⓘ

EDS was not executed - this section will therefore be empty.

6.4 Ligands ⓘ

EDS was not executed - this section will therefore be empty.

6.5 Other polymers ⓘ

EDS was not executed - this section will therefore be empty.