



Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 11:31 AM GMT

PDB ID : 1J2F
Title : X-ray crystal structure of IRF-3 and its functional implications
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Deposited on : 2003-01-04
Resolution : 2.30 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

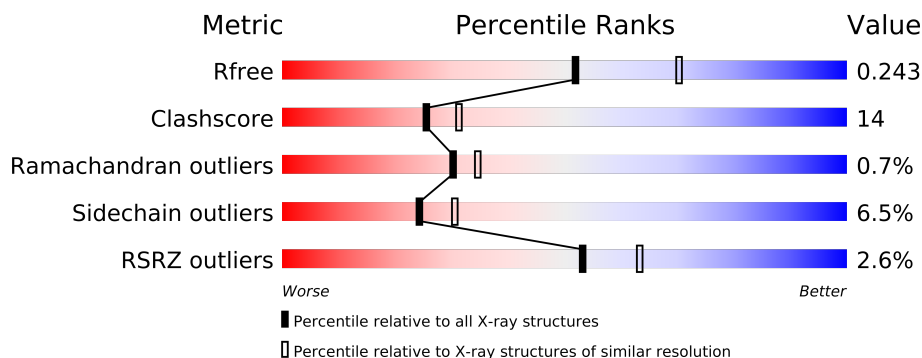
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 2.30 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}	66092	2929 (2.30-2.30)
Clashscore	79885	3679 (2.30-2.30)
Ramachandran outliers	78287	3642 (2.30-2.30)
Sidechain outliers	78261	3641 (2.30-2.30)
RSRZ outliers	66119	2930 (2.30-2.30)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	258	
1	B	258	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3728 atoms, of which 0 are hydrogen and 0 are deuterium.

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 Interferon regulatory factor 3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	226	Total	C	N	O	S	0	0	0
			1757	1126	301	321	9			
1	B	231	Total	C	N	O	S	0	0	0
			1811	1158	310	333	10			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	170	GLY	-	CLONING ARTIFACT	UNP Q14653
A	171	ALA	-	CLONING ARTIFACT	UNP Q14653
A	172	MET	-	CLONING ARTIFACT	UNP Q14653
A	173	GLY	-	CLONING ARTIFACT	UNP Q14653
A	174	SER	-	CLONING ARTIFACT	UNP Q14653
B	170	GLY	-	CLONING ARTIFACT	UNP Q14653
B	171	ALA	-	CLONING ARTIFACT	UNP Q14653
B	172	MET	-	CLONING ARTIFACT	UNP Q14653
B	173	GLY	-	CLONING ARTIFACT	UNP Q14653
B	174	SER	-	CLONING ARTIFACT	UNP Q14653

- Molecule 2 is water.

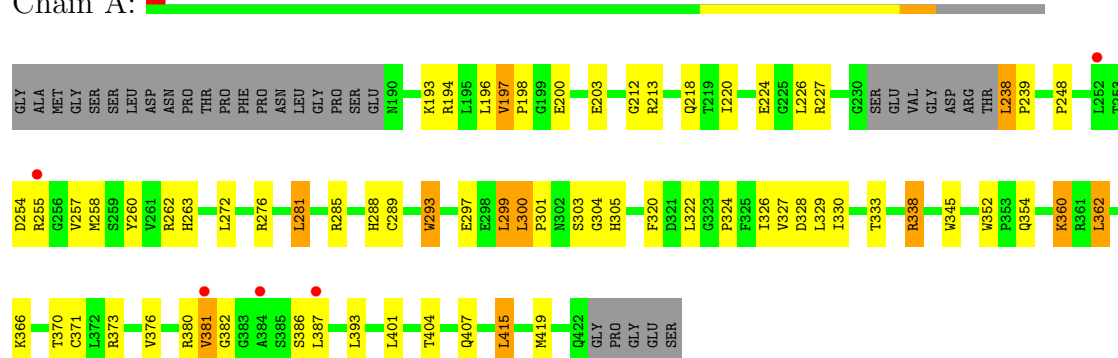
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	75	Total	O	0	0
			75	75		
2	B	85	Total	O	0	0
			85	85		

3 Residue-property plots

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.

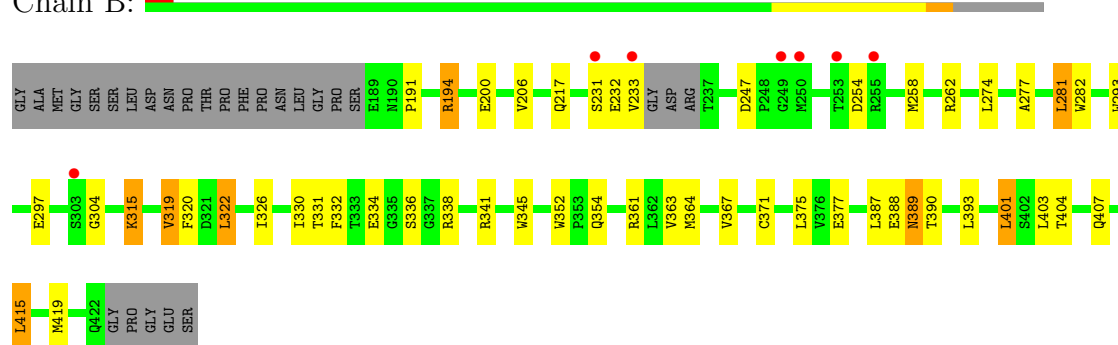
• Molecule 1: Interferon regulatory factor 3

Chain A:



• Molecule 1: Interferon regulatory factor 3

Chain B:



4 Data and refinement statistics

Property	Value	Source
Space group	P 4 21 2	Depositor
Cell constants a, b, c, α , β , γ	134.81Å 134.81Å 69.34Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	62.02 – 2.30 61.66 – 2.00	Depositor EDS
% Data completeness (in resolution range)	(Not available) (62.02-2.30) 94.1 (61.66-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.00 (at 2.00Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.217 , 0.242 0.223 , 0.243	Depositor DCC
R_{free} test set	2677 reflections (9.92%)	DCC
Wilson B-factor (Å ²)	29.7	Xtriage
Anisotropy	0.101	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 39.5	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	1 of 40841 reflections (0.002%)	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	3728	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 13.68% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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	0.34	0/1808	0.58	0/2468
1	B	0.36	0/1863	0.59	0/2540
All	All	0.35	0/3671	0.59	0/5008

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1757	0	1690	55	0
1	B	1811	0	1752	43	0
2	A	75	0	0	4	0
2	B	85	0	0	4	0
All	All	3728	0	3442	95	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 14.

All (95) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:303:SER:O	1:A:305:HIS:N	2.01	0.93

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:326:ILE:HD13	1:B:415:LEU:HD13	1.54	0.89
1:A:338:ARG:HD2	1:A:338:ARG:H	1.38	0.88
1:A:326:ILE:HD13	1:A:415:LEU:HD13	1.59	0.83
1:B:254:ASP:HB3	2:B:42:HOH:O	1.86	0.76
1:B:297:GLU:OE2	1:B:341:ARG:HD3	1.88	0.74
1:A:299:LEU:HD21	1:B:387:LEU:HB2	1.69	0.73
1:A:227:ARG:CZ	1:A:238:LEU:HD13	2.20	0.71
1:A:285:ARG:HH21	1:A:288:HIS:HA	1.57	0.69
1:A:352:TRP:NE1	1:A:354:GLN:HG3	2.09	0.68
1:A:326:ILE:HD12	1:A:419:MET:CE	2.26	0.66
1:A:212:GLY:O	1:B:388:GLU:HG3	1.95	0.65
1:B:331:THR:HG22	1:B:336:SER:HB2	1.78	0.64
1:B:352:TRP:NE1	1:B:354:GLN:HG3	2.13	0.63
1:A:212:GLY:HA2	1:B:387:LEU:HD23	1.80	0.62
1:B:338:ARG:HB3	1:B:338:ARG:HH11	1.65	0.61
1:B:401:LEU:HD22	1:B:403:LEU:HG	1.82	0.60
1:B:277:ALA:HB3	1:B:282:TRP:HE1	1.67	0.59
1:A:387:LEU:O	1:A:404:THR:HA	2.03	0.58
1:B:277:ALA:HB3	1:B:282:TRP:NE1	2.20	0.57
1:A:329:LEU:O	1:A:333:THR:HG23	2.05	0.57
1:A:345:TRP:CZ2	1:A:366:LYS:HD3	2.40	0.57
1:B:326:ILE:CD1	1:B:415:LEU:HD13	2.31	0.56
1:B:281:LEU:HB2	1:B:320:PHE:HB3	1.87	0.56
1:A:218:GLN:NE2	1:A:220:ILE:HD11	2.21	0.55
1:A:381:VAL:HG12	1:A:381:VAL:O	2.06	0.54
1:A:324:PRO:O	1:A:327:VAL:HG22	2.08	0.54
1:B:361:ARG:HB3	1:B:364:MET:CE	2.37	0.54
1:B:330:ILE:O	1:B:334:GLU:HG3	2.07	0.54
1:A:198:PRO:HG2	2:A:133:HOH:O	2.07	0.54
1:A:197:VAL:HG13	1:A:200:GLU:HG3	1.90	0.53
1:A:218:GLN:CD	1:A:220:ILE:HD11	2.29	0.52
1:A:380:ARG:C	1:A:382:GLY:H	2.13	0.52
1:B:322:LEU:HD22	1:B:326:ILE:CD1	2.40	0.52
1:B:232:GLU:HA	2:B:39:HOH:O	2.10	0.51
1:B:293:TRP:CE3	1:B:319:VAL:HG22	2.46	0.51
1:A:276:ARG:HD3	2:A:17:HOH:O	2.11	0.50
1:A:371:CYS:HB3	1:A:393:LEU:O	2.13	0.49
1:A:404:THR:OG1	1:A:407:GLN:HG3	2.13	0.49
1:A:254:ASP:OD2	1:A:257:VAL:HG23	2.13	0.49
1:A:338:ARG:N	1:A:338:ARG:HD2	2.18	0.49
1:A:220:ILE:N	1:A:220:ILE:HD12	2.28	0.49
1:B:206:VAL:O	1:B:217:GLN:HA	2.13	0.49

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:338:ARG:NH1	1:B:338:ARG:HB3	2.28	0.48
1:A:196:LEU:O	1:A:276:ARG:NH2	2.46	0.48
1:B:332:PHE:HA	1:B:336:SER:HB3	1.95	0.48
1:B:403:LEU:HB3	1:B:407:GLN:HG3	1.95	0.48
1:A:415:LEU:HD22	1:A:419:MET:HE2	1.96	0.48
1:B:258:MET:O	1:B:262:ARG:HG3	2.14	0.48
1:A:303:SER:C	1:A:305:HIS:H	2.09	0.48
1:B:361:ARG:HB3	1:B:364:MET:HE3	1.95	0.48
1:A:326:ILE:HD12	1:A:419:MET:HE1	1.94	0.48
1:A:373:ARG:O	1:A:376:VAL:HG22	2.13	0.47
1:A:338:ARG:H	1:A:338:ARG:CD	2.15	0.47
1:B:326:ILE:HD12	1:B:419:MET:CE	2.45	0.47
1:B:231:SER:OG	1:B:233:VAL:HG23	2.15	0.47
1:A:360:LYS:H	1:A:360:LYS:HD3	1.79	0.47
1:A:360:LYS:HE3	2:A:141:HOH:O	2.15	0.46
1:B:388:GLU:HG2	1:B:404:THR:HG22	1.98	0.46
1:B:191:PRO:HG3	1:B:401:LEU:HD11	1.98	0.45
1:B:389:ASN:ND2	1:B:390:THR:H	2.15	0.45
1:A:327:VAL:HG23	1:A:328:ASP:N	2.30	0.44
1:B:247:ASP:OD2	1:B:262:ARG:NH1	2.47	0.44
1:A:285:ARG:HD2	1:A:289:CYS:O	2.17	0.44
1:B:371:CYS:HB3	1:B:393:LEU:O	2.17	0.44
1:A:387:LEU:C	1:A:404:THR:HG22	2.38	0.44
1:A:352:TRP:HE1	1:A:354:GLN:HG3	1.83	0.43
1:B:293:TRP:HA	1:B:345:TRP:O	2.18	0.43
1:A:255:ARG:HG3	2:A:1:HOH:O	2.19	0.43
1:B:274:LEU:CD2	1:B:367:VAL:HG11	2.48	0.43
1:A:226:LEU:C	1:A:226:LEU:HD12	2.39	0.43
1:A:238:LEU:HD12	1:A:238:LEU:N	2.32	0.43
1:B:338:ARG:NH1	1:B:338:ARG:CB	2.82	0.43
1:A:281:LEU:HB2	1:A:320:PHE:HB3	2.01	0.43
1:B:315:LYS:C	1:B:315:LYS:HD2	2.39	0.43
1:A:293:TRP:HA	1:A:345:TRP:O	2.19	0.42
1:B:274:LEU:HA	1:B:282:TRP:O	2.20	0.42
1:A:203:GLU:HB3	1:A:370:THR:OG1	2.19	0.42
1:A:360:LYS:HB2	1:A:362:LEU:HD13	2.01	0.42
1:A:213:ARG:HG3	2:B:19:HOH:O	2.18	0.42
1:B:315:LYS:HD2	1:B:315:LYS:O	2.19	0.42
1:A:248:PRO:HB2	1:A:258:MET:HG2	2.02	0.42
1:A:300:LEU:HD23	1:A:301:PRO:HD2	2.00	0.42
1:A:360:LYS:HE2	1:A:360:LYS:HB2	1.83	0.41
1:B:363:VAL:O	1:B:364:MET:HE2	2.20	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:262:ARG:HH11	1:B:262:ARG:HG2	1.84	0.41
1:A:224:GLU:HG2	1:A:239:PRO:HB2	2.02	0.41
1:A:322:LEU:HG	1:A:326:ILE:CD1	2.50	0.41
1:B:404:THR:OG1	1:B:407:GLN:HG2	2.20	0.41
1:B:194:ARG:HG2	2:B:30:HOH:O	2.21	0.41
1:A:272:LEU:HD13	1:A:272:LEU:C	2.40	0.41
1:A:326:ILE:O	1:A:330:ILE:HG13	2.22	0.40
1:B:194:ARG:O	1:B:200:GLU:HG3	2.20	0.40
1:A:260:TYR:O	1:A:263:HIS:HB2	2.20	0.40
1:A:258:MET:O	1:A:262:ARG:HG2	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone ⓘ

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	222/258 (86%)	212 (96%)	8 (4%)	2 (1%)	25	26
1	B	227/258 (88%)	216 (95%)	10 (4%)	1 (0%)	43	52
All	All	449/516 (87%)	428 (95%)	18 (4%)	3 (1%)	30	34

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	304	GLY
1	A	381	VAL
1	B	304	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	187/217 (86%)	172 (92%)	15 (8%)	17	20
1	B	196/217 (90%)	186 (95%)	10 (5%)	33	43
All	All	383/434 (88%)	358 (94%)	25 (6%)	24	30

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	193	LYS
1	A	194	ARG
1	A	197	VAL
1	A	238	LEU
1	A	281	LEU
1	A	293	TRP
1	A	297	GLU
1	A	299	LEU
1	A	300	LEU
1	A	338	ARG
1	A	360	LYS
1	A	362	LEU
1	A	386	SER
1	A	401	LEU
1	A	415	LEU
1	B	194	ARG
1	B	281	LEU
1	B	315	LYS
1	B	319	VAL
1	B	322	LEU
1	B	375	LEU
1	B	377	GLU
1	B	389	ASN
1	B	401	LEU
1	B	415	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	214	GLN
1	A	284	GLN
1	A	389	ASN
1	B	217	GLN
1	B	284	GLN

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Mol	Chain	Res	Type
1	B	389	ASN
1	B	413	GLN
1	B	422	GLN

5.3.3 RNA ⓘ

There are no RNA chains in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

There are no ligands in this entry.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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	226/258 (87%)	-0.07	5 (2%) 59 69	19, 30, 60, 72	0
1	B	231/258 (89%)	-0.09	7 (3%) 48 58	16, 27, 53, 74	0
All	All	457/516 (88%)	-0.08	12 (2%) 53 63	16, 29, 58, 74	0

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	249	GLY	4.3
1	A	381	VAL	3.6
1	B	250	MET	3.3
1	B	233	VAL	2.9
1	A	255	ARG	2.9
1	B	255	ARG	2.7
1	A	252	LEU	2.7
1	B	303	SER	2.6
1	A	384	ALA	2.6
1	B	253	THR	2.4
1	A	387	LEU	2.2
1	B	231	SER	2.1

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.