



wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 1, 2016 – 09:50 AM GMT

PDB ID : 3JV4
Title : Crystal structure of the dimerization domains p50 and RelB
Authors : Vu, D.; Huang, D.B.; Ghosh, G.
Deposited on : 2009-09-15
Resolution : 3.15 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
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
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) : trunk26865

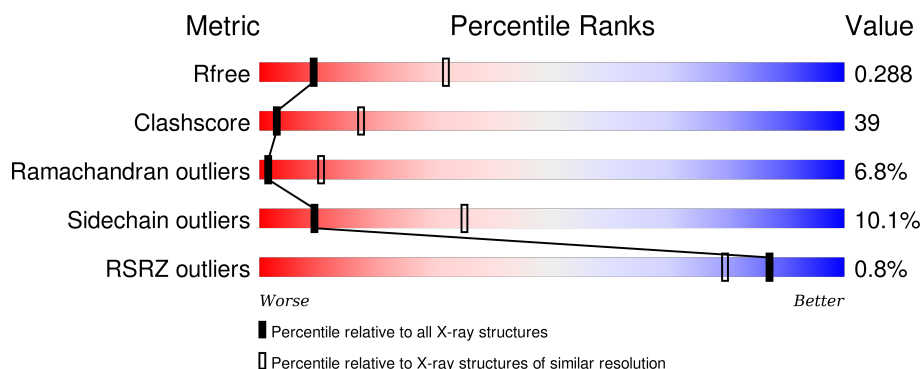
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.15 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}	91344	1112 (3.20-3.12)
Clashscore	102246	1249 (3.20-3.12)
Ramachandran outliers	100387	1222 (3.20-3.12)
Sidechain outliers	100360	1221 (3.20-3.12)
RSRZ outliers	91569	1117 (3.20-3.12)

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	101	<div> <div>2%</div> <div>38%</div> <div>52%</div> <div>9%</div> </div>
1	C	101	<div> <div>%</div> <div>36%</div> <div>55%</div> <div>9%</div> </div>
1	E	101	<div> <div>%</div> <div>38%</div> <div>53%</div> <div>9%</div> </div>
2	B	115	<div> <div>33%</div> <div>56%</div> <div>9%</div> </div>
2	D	115	<div> <div>43%</div> <div>46%</div> <div>10%</div> </div>

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Mol	Chain	Length	Quality of chain
2	F	115	 <div> 37% 53% 7% </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 5185 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 Transcription factor RelB.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	101	Total	C	N	O	S	0	0	0
			800	507	132	157	4			
1	C	101	Total	C	N	O	S	0	0	0
			800	507	132	157	4			
1	E	101	Total	C	N	O	S	0	0	0
			800	507	132	157	4			

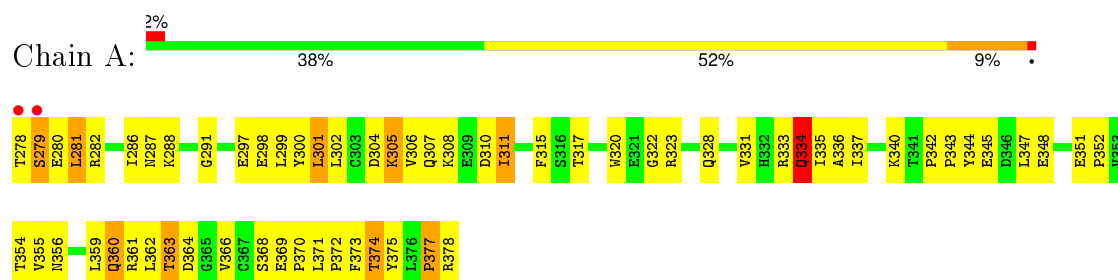
- Molecule 2 is a protein called Nuclear factor NF-kappa-B p105 subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	115	Total	C	N	O	S	0	0	0
			932	594	154	181	3			
2	D	115	Total	C	N	O	S	0	0	0
			942	600	158	181	3			
2	F	112	Total	C	N	O	S	0	0	0
			911	581	150	177	3			

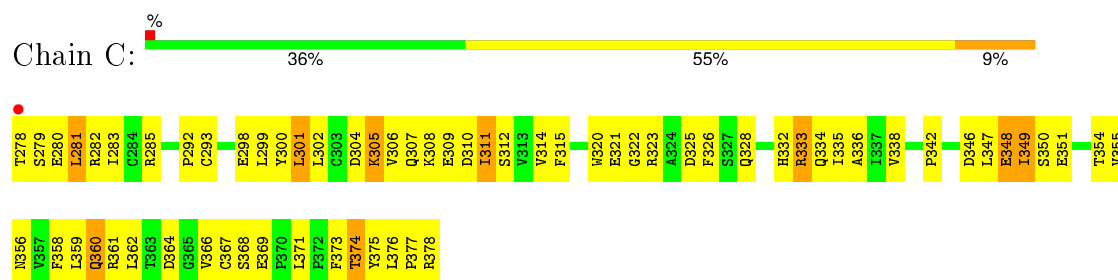
3 Residue-property plots

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.

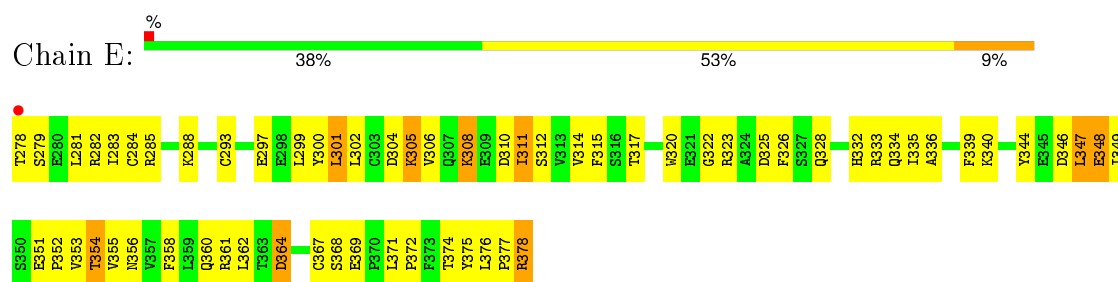
• Molecule 1: Transcription factor RelB



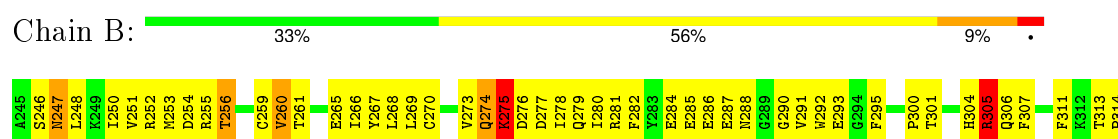
• Molecule 1: Transcription factor RelB



• Molecule 1: Transcription factor RelB



• Molecule 2: Nuclear factor NF-kappa-B p105 subunit





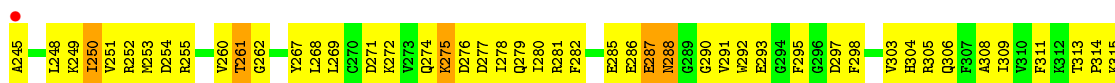
• Molecule 2: Nuclear factor NF-kappa-B p105 subunit

Chain D: 43% 46% 10%



• Molecule 2: Nuclear factor NF-kappa-B p105 subunit

Chain F: 37% 53% 7%



4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	106.17Å 106.17Å 143.78Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	29.76 – 3.15 42.50 – 3.14	Depositor EDS
% Data completeness (in resolution range)	81.6 (29.76-3.15) 90.7 (42.50-3.14)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.04 (at 3.12Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.225 , 0.280 0.231 , 0.288	Depositor DCC
R_{free} test set	754 reflections (4.97%)	DCC
Wilson B-factor (Å ²)	83.3	Xtriage
Anisotropy	0.297	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 45.2	EDS
Estimated twinning fraction	0.035 for -h,-k,l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 16285 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	5185	wwPDB-VP
Average B, all atoms (Å ²)	89.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.34% 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, 0.375 respectively for untwinned datasets, and 0.333, 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	0.39	0/817	0.67	0/1111
1	C	0.43	0/817	0.72	0/1111
1	E	0.35	0/817	0.62	0/1111
2	B	0.49	0/952	0.75	1/1285 (0.1%)
2	D	0.48	0/962	0.75	0/1296
2	F	0.44	0/931	0.66	0/1256
All	All	0.44	0/5296	0.70	1/7170 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	337	LEU	CA-CB-CG	5.38	127.68	115.30

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	800	0	785	68	0
1	C	800	0	785	74	0
1	E	800	0	785	66	0
2	B	932	0	904	101	0
2	D	942	0	926	73	0
2	F	911	0	885	60	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	5185	0	5070	401	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 39.

The worst 5 of 401 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:C:306:VAL:HG11	1:C:311:ILE:HG12	1.41	1.02
1:E:347:LEU:H	1:E:347:LEU:HD12	1.30	0.96
2:B:317:LYS:HG3	2:B:318:ASP:H	1.31	0.94
1:C:293:CYS:HA	1:C:375:TYR:HD2	1.31	0.94
1:C:283:ILE:HG12	1:C:359:LEU:CD1	2.00	0.92

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	99/101 (98%)	79 (80%)	11 (11%)	9 (9%)	1	5
1	C	99/101 (98%)	84 (85%)	10 (10%)	5 (5%)	2	19
1	E	99/101 (98%)	81 (82%)	15 (15%)	3 (3%)	5	34
2	B	113/115 (98%)	82 (73%)	22 (20%)	9 (8%)	1	7
2	D	113/115 (98%)	92 (81%)	11 (10%)	10 (9%)	1	5
2	F	110/115 (96%)	91 (83%)	12 (11%)	7 (6%)	2	12
All	All	633/648 (98%)	509 (80%)	81 (13%)	43 (7%)	1	10

5 of 43 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	279	SER
1	A	377	PRO
2	B	275	LYS
2	B	288	ASN
2	B	305	ARG

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	92/92 (100%)	86 (94%)	6 (6%)	21	60
1	C	92/92 (100%)	83 (90%)	9 (10%)	10	37
1	E	92/92 (100%)	86 (94%)	6 (6%)	21	60
2	B	102/104 (98%)	89 (87%)	13 (13%)	5	24
2	D	104/104 (100%)	88 (85%)	16 (15%)	3	15
2	F	100/104 (96%)	91 (91%)	9 (9%)	12	42
All	All	582/588 (99%)	523 (90%)	59 (10%)	9	35

5 of 59 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	348	GLU
2	D	301	THR
2	F	319	VAL
1	C	374	THR
2	D	250	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 11 such sidechains are listed below:

Mol	Chain	Res	Type
2	B	306	GLN
2	B	320	ASN
1	E	307	GLN
2	B	274	GLN

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Mol	Chain	Res	Type
2	D	306	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 [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	101/101 (100%)	0.07	2 (1%) 68 52	66, 102, 124, 140	0
1	C	101/101 (100%)	-0.22	1 (0%) 84 74	56, 83, 111, 131	0
1	E	101/101 (100%)	-0.04	1 (0%) 84 74	79, 115, 138, 140	0
2	B	115/115 (100%)	-0.17	0 100 100	49, 70, 115, 136	0
2	D	115/115 (100%)	-0.20	0 100 100	51, 71, 116, 140	0
2	F	112/115 (97%)	-0.19	1 (0%) 85 77	57, 88, 115, 135	0
All	All	645/648 (99%)	-0.13	5 (0%) 87 79	49, 87, 131, 140	0

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	278	THR	2.6
1	C	278	THR	2.6
1	A	278	THR	2.3
2	F	245	ALA	2.1
1	A	279	SER	2.1

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 [i](#)

There are no such residues in this entry.