



wwPDB X-ray Structure Validation Summary Report ⓘ

May 17, 2020 – 12:29 pm BST

PDB ID : 3REA
Title : HIV-1 Nef protein in complex with engineered Hck-SH3 domain
Authors : Schulte, A.; Blankenfeldt, W.; Geyer, M.
Deposited on : 2011-04-04
Resolution : 2.00 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

<https://www.wwpdb.org/validation/2017/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)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

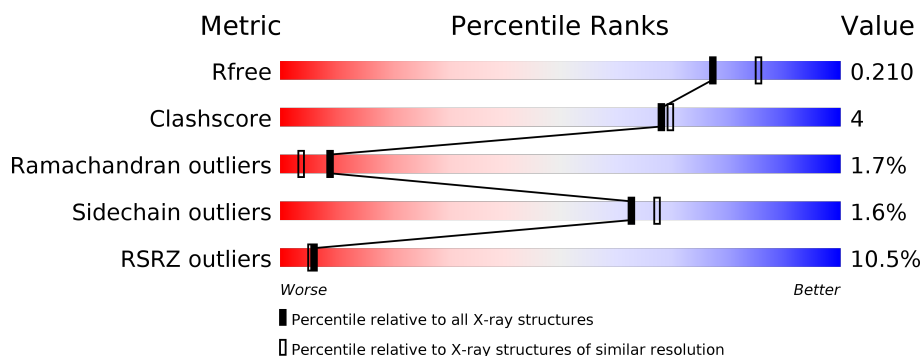
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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}	130704	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

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 respectively. 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	166	<div> <div>8%</div> <div> <div></div> <div>64%</div> <div>10%</div> <div>•</div> <div>23%</div> </div> </div>
1	C	166	<div> <div>11%</div> <div> <div></div> <div>69%</div> <div>7%</div> <div></div> <div>25%</div> </div> </div>
2	B	61	<div> <div>5%</div> <div> <div></div> <div>89%</div> <div>8%</div> <div>• •</div> </div> </div>
2	D	61	<div> <div>7%</div> <div> <div></div> <div>89%</div> <div>10%</div> <div>•</div> </div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3415 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 Protein Nef.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	127	Total	C	N	O	S	0	3	0
			1067	701	181	182	3			
1	C	125	Total	C	N	O	S	0	1	0
			1030	677	176	174	3			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	47	MET	ILE	ENGINEERED MUTATION	UNP P03407
A	48	ALA	THR	ENGINEERED MUTATION	UNP P03407
A	59	SER	CYS	ENGINEERED MUTATION	UNP P03407
A	210	ALA	CYS	ENGINEERED MUTATION	UNP P03407
C	47	MET	ILE	ENGINEERED MUTATION	UNP P03407
C	48	ALA	THR	ENGINEERED MUTATION	UNP P03407
C	59	SER	CYS	ENGINEERED MUTATION	UNP P03407
C	210	ALA	CYS	ENGINEERED MUTATION	UNP P03407

- Molecule 2 is a protein called Tyrosine-protein kinase HCK.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	60	Total	C	N	O	S	0	0	0
			489	312	78	98	1			
2	D	60	Total	C	N	O	S	0	1	0
			498	319	81	96	2			

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	78	MET	-	INITIATING METHIONINE	UNP P08631
B	90	VAL	GLU	ENGINEERED MUTATION	UNP P08631
B	91	SER	ALA	ENGINEERED MUTATION	UNP P08631
B	92	TRP	ILE	ENGINEERED MUTATION	UNP P08631

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Chain	Residue	Modelled	Actual	Comment	Reference
B	93	SER	HIS	ENGINEERED MUTATION	UNP P08631
B	94	PRO	HIS	ENGINEERED MUTATION	UNP P08631
B	95	ASP	GLU	ENGINEERED MUTATION	UNP P08631
D	78	MET	-	INITIATING METHIONINE	UNP P08631
D	90	VAL	GLU	ENGINEERED MUTATION	UNP P08631
D	91	SER	ALA	ENGINEERED MUTATION	UNP P08631
D	92	TRP	ILE	ENGINEERED MUTATION	UNP P08631
D	93	SER	HIS	ENGINEERED MUTATION	UNP P08631
D	94	PRO	HIS	ENGINEERED MUTATION	UNP P08631
D	95	ASP	GLU	ENGINEERED MUTATION	UNP P08631

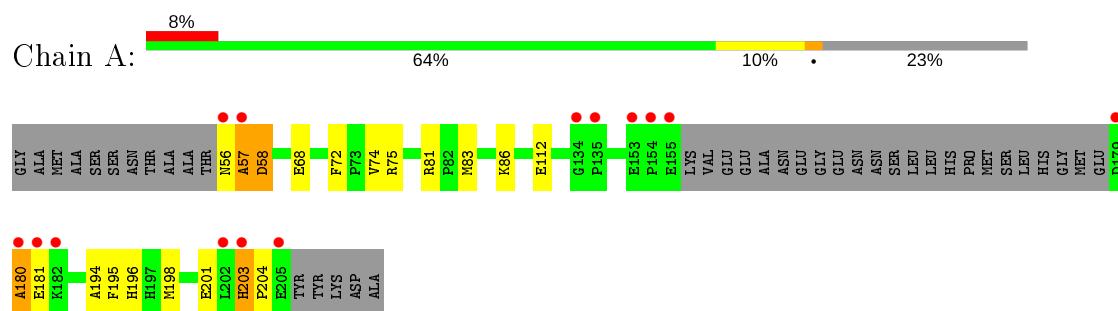
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	129	Total O 129 129	0	0
3	B	52	Total O 52 52	0	0
3	C	89	Total O 89 89	0	0
3	D	61	Total O 61 61	0	0

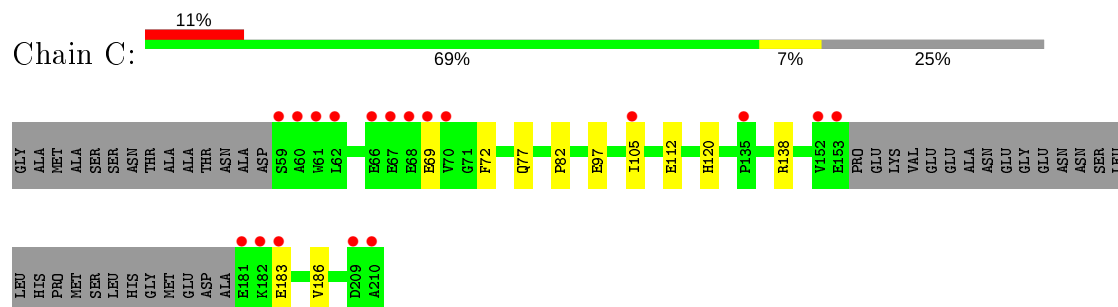
3 Residue-property plots [i](#)

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.

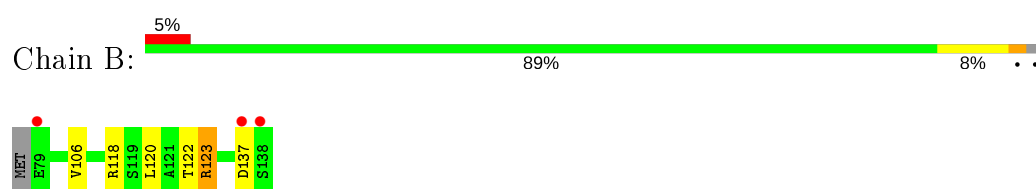
• Molecule 1: Protein Nef



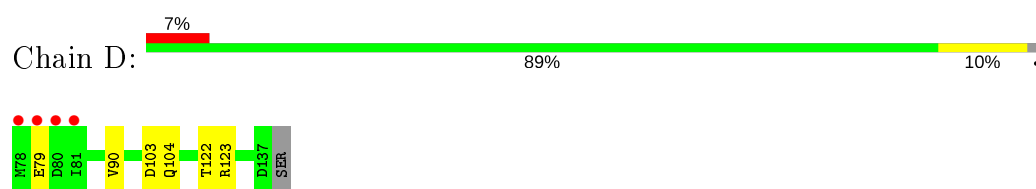
• Molecule 1: Protein Nef



• Molecule 2: Tyrosine-protein kinase HCK



• Molecule 2: Tyrosine-protein kinase HCK



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	65.69 Å 65.69 Å 279.07 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.00 45.82 – 2.00	Depositor EDS
% Data completeness (in resolution range)	100.0 (50.00-2.00) 100.0 (45.82-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.53 (at 2.00 Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.168 , 0.205 0.177 , 0.210	Depositor DCC
R_{free} test set	2148 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	36.0	Xtriage
Anisotropy	0.036	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 60.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	3415	wwPDB-VP
Average B, all atoms (Å ²)	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.29% 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

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	1.29	7/1114 (0.6%)	0.97	2/1516 (0.1%)
1	C	1.15	2/1070 (0.2%)	0.86	0/1455
2	B	1.30	2/501 (0.4%)	1.04	0/680
2	D	1.20	1/513 (0.2%)	0.96	1/696 (0.1%)
All	All	1.23	12/3198 (0.4%)	0.94	3/4347 (0.1%)

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	1

The worst 5 of 12 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	123	ARG	CZ-NH1	9.12	1.45	1.33
2	D	90	VAL	CB-CG1	6.35	1.66	1.52
1	A	194	ALA	CA-CB	6.29	1.65	1.52
1	C	112	GLU	CB-CG	6.13	1.63	1.52
1	C	97	GLU	CG-CD	5.96	1.60	1.51

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	83	MET	CG-SD-CE	8.01	113.01	100.20
2	D	103	ASP	CB-CG-OD1	5.87	123.58	118.30
1	A	81	ARG	NE-CZ-NH2	-5.32	117.64	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	58	ASP	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	1067	0	1019	14	0
1	C	1030	0	974	4	0
2	B	489	0	462	3	0
2	D	498	0	479	2	0
3	A	129	0	0	4	1
3	B	52	0	0	0	0
3	C	89	0	0	1	0
3	D	61	0	0	1	0
All	All	3415	0	2934	23	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 4.

The worst 5 of 23 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:196:HIS:HB3	1:A:198:MET:HE2	1.29	1.10
1:A:196:HIS:HB3	1:A:198:MET:CE	1.91	1.01
1:A:180:ALA:HB3	1:A:181:GLU:HA	1.51	0.91
1:A:75[A]:ARG:NH1	3:A:364:HOH:O	2.18	0.77
1:A:180:ALA:CB	1:A:181:GLU:HA	2.25	0.66

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)
3:A:395:HOH:O	3:A:395:HOH:O[7_555]	1.66	0.54

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	126/166 (76%)	122 (97%)	0	4 (3%)	4	1
1	C	122/166 (74%)	120 (98%)	1 (1%)	1 (1%)	19	13
2	B	58/61 (95%)	55 (95%)	2 (3%)	1 (2%)	9	4
2	D	59/61 (97%)	58 (98%)	1 (2%)	0	100	100
All	All	365/454 (80%)	355 (97%)	4 (1%)	6 (2%)	9	4

5 of 6 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	57	ALA
1	A	203	HIS
1	A	204	PRO
1	A	180	ALA
2	B	137	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	109/141 (77%)	106 (97%)	3 (3%)	43	44
1	C	102/141 (72%)	101 (99%)	1 (1%)	76	81
2	B	53/54 (98%)	53 (100%)	0	100	100
2	D	54/54 (100%)	53 (98%)	1 (2%)	57	61
All	All	318/390 (82%)	313 (98%)	5 (2%)	62	67

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	58	ASP
1	A	72	PHE
1	A	203	HIS
1	C	72	PHE
2	D	79	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	104	GLN
1	C	77	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	127/166 (76%)	0.46	14 (11%) 5 5	36, 44, 95, 111	0
1	C	125/166 (75%)	0.74	18 (14%) 2 2	35, 54, 127, 141	0
2	B	60/61 (98%)	0.08	3 (5%) 28 28	37, 48, 86, 102	0
2	D	60/61 (98%)	0.18	4 (6%) 17 17	36, 48, 80, 111	0
All	All	372/454 (81%)	0.45	39 (10%) 6 5	35, 48, 102, 141	0

The worst 5 of 39 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	180	ALA	9.7
1	C	210	ALA	8.1
1	C	62	LEU	7.7
2	D	78	MET	7.5
1	A	154	PRO	6.8

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.