

wwPDB X-ray Structure Validation Summary Report (i)

Sep 26, 2023 – 05:32 PM EDT

PDB ID : 6CDX

Title: High-resolution crystal structure of fluoropropylated cystine knot, binding to

alpha-5 beta-6 integrin

Authors: Kimura, R.; Nix, J.; Bongura, C.; Chakraborti, S.; Gambhir, S.; Filipp, F.V.

Deposited on : 2018-02-09

Resolution : 1.00 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 (i) symbol.

The types of validation reports are described at http://www.wwpdb.org/validation/2017/FAQs#types.

The following versions of software and data (see references (1)) were used in the production of this report:

 $Mol Probity \quad : \quad 4.02b\text{--}467$

Mogul: 1.8.5 (274361), CSD as541be (2020)

Xtriage (Phenix) : 1.13 EDS : 2.35.1

Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)

 $Refmac \quad : \quad 5.8.0158$

CCP4 : 7.0.044 (Gargrove)

Ideal geometry (proteins) : Engh & Huber (2001) Ideal geometry (DNA, RNA) : Parkinson et al. (1996)

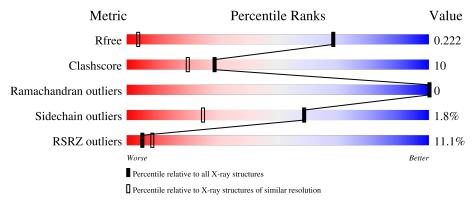
Validation Pipeline (wwPDB-VP) : 2.35.1

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: X-RAY DIFFRACTION

The reported resolution of this entry is 1.00 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	Similar resolution
Metric	$(\# \mathrm{Entries})$	$(\# ext{Entries}, ext{ resolution range}(\mathring{A}))$
R_{free}	130704	1050 (1.06-0.94)
Clashscore	141614	1117 (1.06-0.94)
Ramachandran outliers	138981	1043 (1.06-0.94)
Sidechain outliers	138945	1045 (1.06-0.94)
RSRZ outliers	127900	1023 (1.06-0.94)

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 of 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 <=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	37	89%	11%		
1	В	37	19% 78%	16% • •		



2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 589 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 cystine knot (fluoropropylated).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf	Trace
1	A	37	Total C F N O S 271 160 1 54 50 6	0	1	0
1	В	36	Total C N O S 266 157 54 49 6	0	1	0

• Molecule 2 is water.

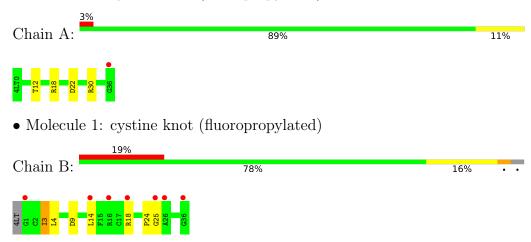
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	22	Total O 22 22	0	0
2	В	30	Total O 30 30	0	0



3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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: cystine knot (fluoropropylated)





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants	16.95Å 44.25Å 66.41Å	Depositor
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor
Resolution (Å)	36.82 - 1.00	Depositor
rtesolution (A)	22.12 - 1.00	EDS
% Data completeness	96.7 (36.82-1.00)	Depositor
(in resolution range)	96.7 (22.12-1.00)	EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	1.67 (at 1.00Å)	Xtriage
Refinement program	REFMAC 5.8.0123	Depositor
D D.	0.203 , 0.226	Depositor
R, R_{free}	0.203 , 0.222	DCC
R_{free} test set	1341 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å ²)	6.5	Xtriage
Anisotropy	0.092	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.37, 35.6	EDS
L-test for twinning ²	$ < L > = 0.48, < L^2> = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	589	wwPDB-VP
Average B, all atoms (Å ²)	9.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 17.56% of the height of the origin peak. No significant pseudotranslation is detected.

²Theoretical values of <|L|>, $<L^2>$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.



¹Intensities estimated from amplitudes.

5 Model quality (i)

5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: 4LT

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		nd lengths	Bond angles	
MIOI	Chain	RMSZ	# Z > 5	RMSZ	# Z >5
1	A	1.31	$2/267 \ (0.7\%)$	1.26	2/354~(0.6%)
1	В	1.25	$1/267 \ (0.4\%)$	1.42	4/354 (1.1%)
All	All	1.28	3/534 (0.6%)	1.35	6/708 (0.8%)

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 maintenain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	В	1	0

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$\mathbf{Observed}(\mathbf{\mathring{A}})$	Ideal(A)
1	A	18	ARG	NE-CZ	-5.76	1.25	1.33
1	A	18	ARG	CZ-NH2	-5.24	1.26	1.33
1	В	25	GLY	C-N	-5.20	1.22	1.34

The worst 5 of 6 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$Observed(^o)$	$Ideal(^{o})$
1	В	18	ARG	NE-CZ-NH2	-11.58	114.51	120.30
1	A	30	ARG	NE-CZ-NH2	6.81	123.71	120.30
1	В	14	LEU	CB-CG-CD2	6.75	122.47	111.00
1	A	22	ASP	CB-CG-OD1	6.67	124.31	118.30
1	В	25	GLY	C-N-CA	5.84	136.30	121.70

All (1) chirality outliers are listed below:



Mol	Chain	Res	Type	Atom
1	В	3	ILE	СВ

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	271	0	252	2	0
1	В	266	0	253	8	0
2	A	22	0	0	2	1
2	В	30	0	0	0	3
All	All	589	0	505	10	3

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 10.

The worst 5 of 10 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$\begin{array}{c} {\rm Interatomic} \\ {\rm distance} \ ({\rm \AA}) \end{array}$	Clash overlap (Å)	
1:B:3:ILE:HD12	1:B:4:LEU:HG	1.68	0.75	
1:B:3:ILE:CD1	1:B:4:LEU:HG	2.17	0.74	
1:A:12:THR:HG22	2:A:103:HOH:O	1.97	0.64	
1:A:12:THR:HG21	2:A:111:HOH:O	2.03	0.59	
1:B:3:ILE:HD12	1:B:4:LEU:N	2.21	0.56	

All (3) 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	$\begin{array}{c} {\rm Interatomic} \\ {\rm distance} \ ({\rm \AA}) \end{array}$	Clash overlap (Å)
2:A:107:HOH:O	2:B:112:HOH:O[2_355]	1.46	0.74
2:B:108:HOH:O	2:B:115:HOH:O[1_455]	1.48	0.72
2:B:115:HOH:O	2:B:128:HOH:O[1_655]	2.16	0.04



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	Percer	ntiles
1	A	35/37~(95%)	33 (94%)	2 (6%)	0	100	100
1	В	$34/37 \ (92\%)$	32 (94%)	2 (6%)	0	100	100
All	All	69/74 (93%)	65 (94%)	4 (6%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains (i)

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	Perce	ntiles
1	A	28/28 (100%)	28 (100%)	0	100	100
1	В	28/28 (100%)	27 (96%)	1 (4%)	35	6
All	All	56/56 (100%)	55 (98%)	1 (2%)	59	25

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	В	3	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 monosaccharides 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, 95^{th} 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	$OWAB(\AA^2)$	Q < 0.9
1	A	$36/37 \ (97\%)$	0.96	1 (2%) 53 44	5, 6, 13, 17	0
1	В	$36/37 \ (97\%)$	1.28	7 (19%) 1 4	4, 8, 16, 22	0
All	All	72/74~(97%)	1.12	8 (11%) 5 8	4, 8, 13, 22	0

The worst 5 of 8 RSRZ outliers are listed below:

Mol	Chain	Res	\mathbf{Type}	RSRZ
1	В	1	GLY	4.7
1	В	26	ALA	3.9
1	В	18	ARG	3.6
1	В	36[A]	GLY	3.2
1	В	14	LEU	2.8

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 monosaccharides 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.

