



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 07:22 am GMT

PDB ID : 5K6H
Title : Crystal structure of prefusion-stabilized RSV F single-chain 9-10 DS-Cav1 A149C-Y458C variant.
Authors : Joyce, M.G.; Zhang, B.; Mascola, J.R.; Kwong, P.D.
Deposited on : 2016-05-24
Resolution : 2.65 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
Xtriage (Phenix) : 1.9-1692
EDS : trunk28620
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
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) : recalc28949

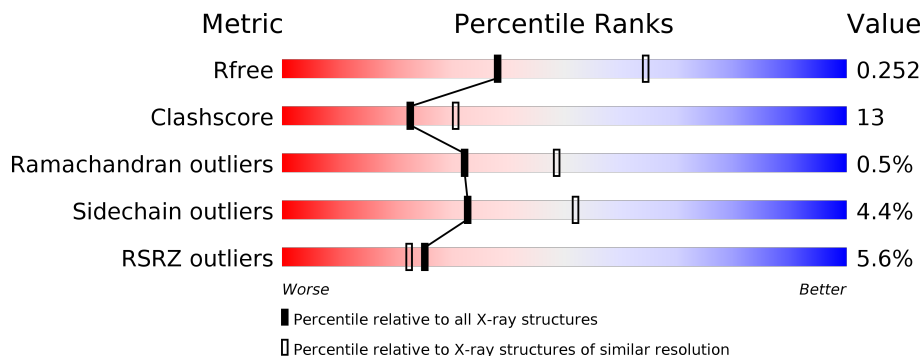
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.65 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}	100719	1044 (2.66-2.62)
Clashscore	112137	1092 (2.66-2.62)
Ramachandran outliers	110173	1077 (2.66-2.62)
Sidechain outliers	110143	1077 (2.66-2.62)
RSRZ outliers	101464	1047 (2.66-2.62)

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	F	445	<div> <div>6%</div> <div>80%</div> <div>18%</div> <div>.</div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3492 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 Fusion glycoprotein F0.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	F	445	Total	C	N	O	S	0	1	0
			3445	2164	576	680	25			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
F	102	ALA	PRO	engineered mutation	UNP P03420
F	143	GLY	-	linker	UNP P03420
F	144	SER	-	linker	UNP P03420
F	149	CYS	ALA	engineered mutation	UNP P03420
F	155	CYS	SER	engineered mutation	UNP P03420
F	190	PHE	SER	engineered mutation	UNP P03420
F	207	LEU	VAL	engineered mutation	UNP P03420
F	290	CYS	SER	engineered mutation	UNP P03420
F	373	ARG	LEU	engineered mutation	UNP P03420
F	379	VAL	ILE	engineered mutation	UNP P03420
F	447	VAL	MET	engineered mutation	UNP P03420
F	458	CYS	TYR	engineered mutation	UNP P03420

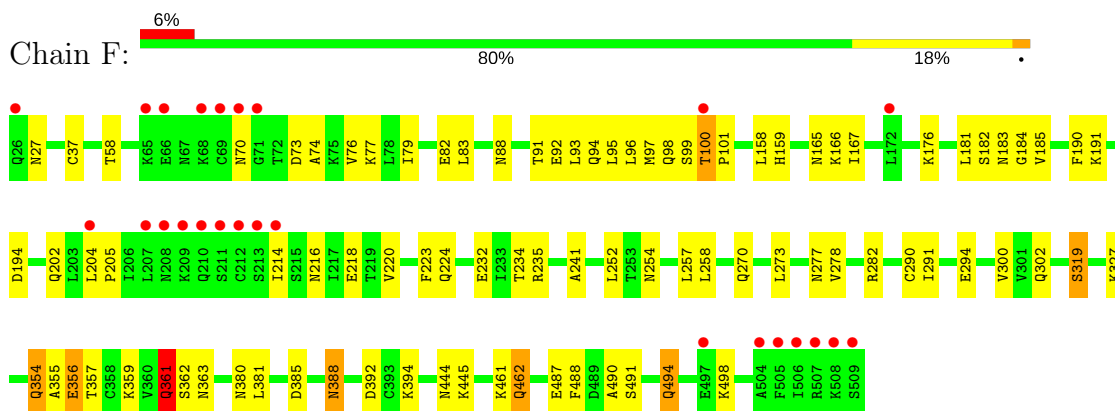
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	F	47	Total	O	0	0
			47	47		

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 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: Fusion glycoprotein F0



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 3 2	Depositor
Cell constants a, b, c, α , β , γ	169.08Å 169.08Å 169.08Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	37.81 – 2.65 37.81 – 2.65	Depositor EDS
% Data completeness (in resolution range)	79.4 (37.81-2.65) 75.4 (37.81-2.65)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.40 (at 2.65Å)	Xtriage
Refinement program	PHENIX (1.10_2155: ???)	Depositor
R, R_{free}	0.222 , 0.252 0.223 , 0.252	Depositor DCC
R_{free} test set	966 reflections (4.94%)	DCC
Wilson B-factor (Å ²)	49.6	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 47.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	3492	wwPDB-VP
Average B, all atoms (Å ²)	70.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.01% 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 [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	F	0.35	0/3497	0.50	0/4737

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	F	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	F	361	GLN	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	F	3445	0	3487	90	0
2	F	47	0	0	5	0
All	All	3492	0	3487	90	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 13.

All (90) 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:F:392:ASP:OD1	1:F:491:SER:HB2	1.54	1.07
1:F:183:ASN:CB	1:F:185:VAL:HG23	1.84	1.06
1:F:99:SER:C	1:F:101:PRO:HD3	1.83	0.99
1:F:183:ASN:HB2	1:F:185:VAL:HG23	1.45	0.98
1:F:94:GLN:O	1:F:97:MET:HG2	1.68	0.93
1:F:99:SER:O	1:F:101:PRO:HD3	1.68	0.92
1:F:73:ASP:OD2	1:F:214:ILE:HG23	1.75	0.85
1:F:183:ASN:N	1:F:184:GLY:HA2	1.92	0.84
1:F:181:LEU:O	1:F:182:SER:OG	1.96	0.83
1:F:183:ASN:HB2	1:F:185:VAL:N	1.96	0.81
1:F:181:LEU:HD13	1:F:185:VAL:O	1.80	0.80
1:F:91:THR:O	1:F:95:LEU:HD13	1.81	0.80
1:F:277:ASN:OD1	1:F:361:GLN:HG2	1.83	0.78
1:F:99:SER:O	1:F:101:PRO:CD	2.31	0.78
1:F:73:ASP:HB3	1:F:76:VAL:HG23	1.65	0.78
1:F:82:GLU:HB3	1:F:223:PHE:HE2	1.50	0.75
1:F:181:LEU:N	1:F:181:LEU:HD12	2.03	0.73
1:F:82:GLU:HB3	1:F:223:PHE:CE2	2.23	0.73
1:F:183:ASN:HD22	1:F:185:VAL:HG21	1.58	0.68
1:F:491:SER:OG	1:F:494:GLN:HB3	1.95	0.67
1:F:183:ASN:HD22	1:F:185:VAL:CG2	2.08	0.66
1:F:79:ILE:HG22	1:F:83:LEU:CD1	2.26	0.66
1:F:183:ASN:HB2	1:F:185:VAL:H	1.59	0.66
1:F:158:LEU:HD21	1:F:167:ILE:HD11	1.77	0.66
1:F:79:ILE:HG22	1:F:83:LEU:HG	1.78	0.65
1:F:361:GLN:HA	1:F:361:GLN:HE21	1.64	0.63
1:F:73:ASP:O	1:F:74:ALA:C	2.32	0.62
1:F:37:CYS:SG	1:F:319:SER:HB3	2.39	0.62
1:F:362:SER:O	1:F:363:ASN:HB3	1.99	0.62
1:F:79:ILE:HG22	1:F:83:LEU:HD11	1.82	0.62
1:F:73:ASP:OD2	1:F:214:ILE:CG2	2.45	0.62
1:F:101:PRO:HB2	1:F:241:ALA:HB1	1.81	0.61
1:F:235:ARG:NH1	2:F:603:HOH:O	2.33	0.60
1:F:183:ASN:HB3	1:F:185:VAL:HG23	1.78	0.60
1:F:355:ALA:O	1:F:356:GLU:HB2	2.00	0.60
1:F:79:ILE:HG22	1:F:83:LEU:CG	2.32	0.59
1:F:100:THR:O	1:F:100:THR:HG22	2.02	0.59
1:F:159:HIS:CE1	1:F:291:ILE:CG2	2.86	0.58

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:392:ASP:OD1	1:F:491:SER:CB	2.42	0.58
1:F:158:LEU:CD2	1:F:167:ILE:HD11	2.33	0.58
1:F:183:ASN:ND2	1:F:185:VAL:CG2	2.68	0.56
1:F:159:HIS:CE1	1:F:291:ILE:HG22	2.42	0.55
1:F:392:ASP:CG	1:F:491:SER:HB2	2.24	0.55
1:F:273:LEU:O	1:F:277:ASN:ND2	2.38	0.55
1:F:101:PRO:CB	1:F:241:ALA:HB1	2.38	0.54
1:F:181:LEU:C	1:F:182:SER:HG	2.05	0.53
1:F:191:LYS:HE3	1:F:194:ASP:OD2	2.10	0.52
1:F:158:LEU:CD2	1:F:167:ILE:CD1	2.89	0.51
1:F:97:MET:HG3	1:F:98:GLN:N	2.25	0.51
1:F:361:GLN:O	1:F:362:SER:C	2.48	0.51
1:F:494:GLN:O	1:F:498:LYS:HB2	2.11	0.51
1:F:181:LEU:N	1:F:181:LEU:CD1	2.73	0.50
1:F:258:LEU:HD21	1:F:278:VAL:HG21	1.94	0.50
1:F:100:THR:N	1:F:101:PRO:HD3	2.28	0.49
1:F:99:SER:C	1:F:101:PRO:CD	2.69	0.48
1:F:270:GLN:NE2	2:F:609:HOH:O	2.46	0.48
1:F:354:GLN:HB2	2:F:628:HOH:O	2.14	0.48
1:F:392:ASP:CG	1:F:491:SER:CB	2.82	0.48
1:F:91:THR:O	1:F:95:LEU:CD1	2.57	0.47
1:F:79:ILE:HD11	1:F:220:VAL:HG22	1.96	0.47
1:F:183:ASN:ND2	1:F:185:VAL:HG23	2.29	0.47
1:F:257:LEU:HD23	1:F:278:VAL:HG13	1.97	0.47
1:F:183:ASN:CG	1:F:185:VAL:HG23	2.34	0.46
1:F:158:LEU:HD22	1:F:167:ILE:HD12	1.98	0.46
1:F:490:ALA:HB1	1:F:494:GLN:HG3	1.97	0.45
1:F:82:GLU:CD	1:F:224:GLN:HG3	2.37	0.45
1:F:74:ALA:O	1:F:77:LYS:N	2.49	0.45
1:F:204:LEU:N	1:F:205:PRO:HD2	2.32	0.44
1:F:290:CYS:SG	1:F:300:VAL:HG23	2.57	0.44
1:F:93:LEU:HD11	1:F:234:THR:HG22	2.01	0.43
1:F:490:ALA:HB1	1:F:494:GLN:CG	2.48	0.43
1:F:27:ASN:N	2:F:606:HOH:O	2.52	0.43
1:F:462:GLN:NE2	2:F:610:HOH:O	2.51	0.42
1:F:216:ASN:ND2	1:F:218:GLU:OE1	2.45	0.42
1:F:354:GLN:HB3	1:F:357:THR:HG23	2.02	0.42
1:F:158:LEU:HD22	1:F:167:ILE:CD1	2.49	0.42
1:F:252:LEU:O	1:F:282:ARG:NH1	2.39	0.42
1:F:258:LEU:HD21	1:F:278:VAL:CG2	2.51	0.41
1:F:79:ILE:O	1:F:83:LEU:HG	2.20	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:444:ASN:O	1:F:461:LYS:NZ	2.54	0.41
1:F:381:LEU:HB3	1:F:388:ASN:OD1	2.20	0.41
1:F:165:ASN:ND2	1:F:294:GLU:OE1	2.53	0.41
1:F:385:ASP:O	1:F:388:ASN:OD1	2.38	0.41
1:F:99:SER:O	1:F:101:PRO:HD2	2.15	0.41
1:F:88:ASN:O	1:F:92:GLU:HG3	2.20	0.41
1:F:232:GLU:OE1	1:F:235:ARG:NH2	2.55	0.40
1:F:94:GLN:O	1:F:97:MET:CG	2.54	0.40
1:F:327:LYS:HD2	1:F:327:LYS:H	1.86	0.40
1:F:58:THR:N	1:F:190:PHE:O	2.51	0.40
1:F:487:GLU:OE1	1:F:494:GLN:NE2	2.53	0.40

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	F	444/445 (100%)	428 (96%)	14 (3%)	2 (0%)	32 48

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	F	70	ASN
1	F	100	THR

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	Percentiles	
1	F	408/407 (100%)	390 (96%)	18 (4%)	33	52

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	F	96	LEU
1	F	166	LYS
1	F	176	LYS
1	F	202	GLN
1	F	254	ASN
1	F	302	GLN
1	F	319	SER
1	F	354	GLN
1	F	356	GLU
1	F	359	LYS
1	F	361	GLN
1	F	380	ASN
1	F	388	ASN
1	F	394	LYS
1	F	445	LYS
1	F	462	GLN
1	F	488	PHE
1	F	494	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	F	361	GLN

5.3.3 RNA ⓘ

There are no RNA molecules 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 [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

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	F	445/445 (100%)	0.10	25 (5%)	25 22	30, 58, 144, 264	0

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	210	GLN	8.7
1	F	211	SER	7.4
1	F	26	GLN	6.4
1	F	214	ILE	5.3
1	F	208	ASN	5.0
1	F	69	CYS	4.8
1	F	207	LEU	4.7
1	F	509	SER	4.7
1	F	507	ARG	4.7
1	F	66	GLU	4.6
1	F	204	LEU	4.3
1	F	212	CYS	4.0
1	F	70	ASN	3.9
1	F	68	LYS	3.9
1	F	71	GLY	3.6
1	F	209	LYS	3.5
1	F	213	SER	2.9
1	F	506	ILE	2.8
1	F	505	PHE	2.8
1	F	508	LYS	2.6
1	F	504	ALA	2.4
1	F	172	LEU	2.2
1	F	65	LYS	2.2
1	F	100	THR	2.1
1	F	497	GLU	2.0

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.