

wwPDB X-ray Structure Validation Summary Report (i)

Oct 10, 2023 – 06:17 AM EDT

PDB ID	:	7KFA
Title	:	PCSK9 in complex with PCSK9i a 13mer cyclic peptide LDLR disruptor
Authors	:	Chopra, R.; Xu, M.; Spraggon, G.
Deposited on	:	2020-10-13
Resolution	:	2.45 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:

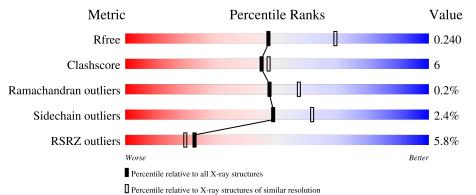
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.35.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (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: $X\text{-}RAY \, DIFFRACTION$

The reported resolution of this entry is 2.45 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	$egin{array}{c} { m Whole \ archive} \ (\#{ m Entries}) \end{array}$	${f Similar\ resolution}\ (\#{ m Entries,\ resolution\ range}({ m \AA}))$
R_{free}	130704	1544 (2.48-2.44)
Clashscore	141614	1613 (2.48-2.44)
Ramachandran outliers	138981	1598 (2.48-2.44)
Sidechain outliers	138945	1598 (2.48-2.44)
RSRZ outliers	127900	1523 (2.48-2.44)

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	А	124	^{2%} 67%	7%	26%
2	В	546	73%	12%	• 14%
3	D	13	62%	31%	8%



2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 4463 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 Proprotein convertase subtilisin/kexin type 9 Propeptide.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	А	92	Total 748	C 479	N 136	O 131	${ m S} { m 2}$	0	1	0

• Molecule 2 is a protein called Proprotein convertase subtilisin/kexin type 9.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace	
2	В	471	Total 3497	C 2154	N 645	O 667	S 31	0	0	0

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
В	474	ILE	VAL	variant	UNP Q8NBP7
В	670	GLU	GLY	variant	UNP Q8NBP7
В	693	HIS	-	expression tag	UNP Q8NBP7
В	694	HIS	-	expression tag	UNP Q8NBP7
В	695	HIS	-	expression tag	UNP Q8NBP7
В	696	HIS	-	expression tag	UNP Q8NBP7
В	697	HIS	-	expression tag	UNP Q8NBP7
В	698	HIS	-	expression tag	UNP Q8NBP7

• Molecule 3 is a protein called 1-[2,6,10.14-TETRAMETHYL-HEXADECAN-16-YL]-2-[2,10 ,14-TRIMETHYLHEXADECAN-16-YL]GLYCEROL.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
3	D	13	Total 118	$\begin{array}{c} \mathrm{C} \\ 85 \end{array}$	N 14	O 18	S 1	0	0	0

• Molecule 4 is CALCIUM ION (three-letter code: CA) (formula: Ca).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	В	1	Total Ca 1 1	0	0

• Molecule 5 is water.

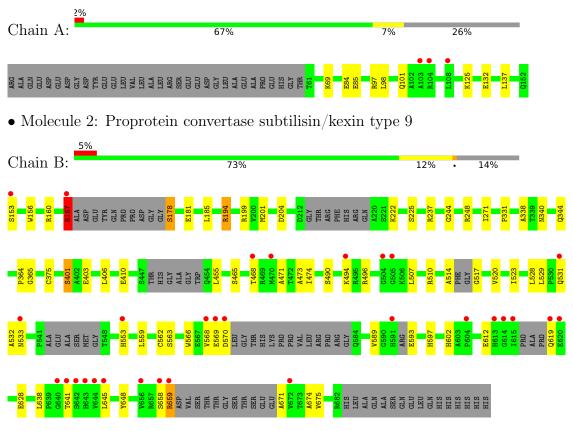
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	А	27	Total O 27 27	0	0
5	В	70	Total O 70 70	0	0
5	D	2	Total O 2 2	0	0



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: Proprotein convertase subtilisin/kexin type 9 Propeptide



• Molecule 3: 1-[2,6,10.14-TETRAMETHYL-HEXADECAN-16-YL]-2-[2,10,14-TRIMETHYLHE XADECAN-16-YL]GLYCEROL





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants	62.42Å 70.34Å 150.62Å	Depositor
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor
Resolution (Å)	57.67 - 2.45	Depositor
Resolution (A)	57.67 - 2.45	EDS
% Data completeness	93.0 (57.67-2.45)	Depositor
(in resolution range)	93.0(57.67-2.45)	EDS
R _{merge}	0.08	Depositor
R _{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	$2.30 (at 2.45 \text{\AA})$	Xtriage
Refinement program	PHENIX 1.17.1_3660	Depositor
D D.	0.189 , 0.242	Depositor
R, R_{free}	0.189 , 0.240	DCC
R_{free} test set	1193 reflections (5.11%)	wwPDB-VP
Wilson B-factor $(Å^2)$	45.6	Xtriage
Anisotropy	0.134	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.36 , 50.2	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	0.95	EDS
Total number of atoms	4463	wwPDB-VP
Average B, all atoms $(Å^2)$	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 4.72% of the height of the origin peak. No significant pseudotranslation is detected.

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.



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: NEH, WCM, ACE, CA

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	Bo	nd lengths	Bond angles		
	Unam	RMSZ	# Z > 5	RMSZ	# Z > 5	
1	А	0.27	0/768	0.47	0/1037	
2	В	0.35	0/3554	0.53	1/4820~(0.0%)	
3	D	1.35	1/68~(1.5%)	1.26	0/93	
All	All	0.37	1/4390~(0.0%)	0.54	1/5950~(0.0%)	

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	9	GLU	CD-OE1	7.99	1.34	1.25

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
2	В	167	ARG	NE-CZ-NH2	5.67	123.13	120.30

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	А	748	0	763	8	0
2	В	3497	0	3416	47	3
3	D	118	0	68	2	0



	0	Non-H	1 0	H(added)	Clashes	Symm-Clashes
	Onam	11011-11	model	II(auucu)	Clashes	Bymm-Clashes
4	В	1	0	0	0	0
5	А	27	0	0	4	0
5	В	70	0	0	4	0
5	D	2	0	0	0	0
All	All	4463	0	4247	56	3

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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 6.

The worst 5 of 56 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)
2:B:641:THR:HG1	2:B:671:ALA:N	1.72	0.88
2:B:194:ARG:O	2:B:237:ARG:NH2	2.15	0.80
2:B:628:GLU:N	2:B:628:GLU:OE1	2.22	0.71
2:B:490:SER:HB2	2:B:520:VAL:HG12	1.77	0.66
2:B:201:MET:HE2	2:B:248:ARG:HH21	1.61	0.65

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	Interatomic distance (Å)	Clash overlap (Å)
2:B:167:ARG:NH1	2:B:612:GLU:CD[4_477]	1.97	0.23
2:B:153:SER:OG	2:B:553:HIS:NE2[1_455]	2.14	0.06
2:B:167:ARG:NH1	2:B:612:GLU:CG[4_477]	2.15	0.05

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	А	91/124~(73%)	89(98%)	2(2%)	0	100 100



	Chain	Analysed	Favoured	Allowed	Outliers	Percentil	es
2	В	451/546~(83%)	439 (97%)	11 (2%)	1 (0%)	47 57]
3	D	7/13~(54%)	7 (100%)	0	0	100 10	0
All	All	549/683~(80%)	535~(97%)	13 (2%)	1 (0%)	47 57	

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All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	В	331	PRO

5.3.2 Protein sidechains (i)

In the following table, the Percentiles column shows the percent side chain 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 side chain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	А	80/104~(77%)	80 (100%)	0	100 100
2	В	379/437~(87%)	369~(97%)	10 (3%)	46 58
3	D	8/8 (100%)	7~(88%)	1 (12%)	4 3
All	All	467/549~(85%)	456~(98%)	11 (2%)	49 61

5 of 11 residues with a non-rotameric side chain are listed below:

Mol	Chain	Res	Type
2	В	570	ASP
2	В	593	GLU
3	D	6	THR
2	В	659	ARG
2	В	204	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
2	В	344	GLN
2	В	513	ASN
2	В	533	ASN



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Mol	Chain	Res	Type
2	В	537	HIS
2	В	643	HIS

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

2 non-standard protein/DNA/RNA residues are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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| > 2 is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mal	Mol Type Chain	Chain	Chain	Chain	Chain	Chain	Chain	Chain	Chain	Dec	Link	Bo	ond leng	\mathbf{ths}	В	ond ang	les
IVIOI		Res	LIIIK	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2								
3	WCM	D	8	3	23,24,25	0.68	0	30,32,34	1.53	4 (13%)							
3	WCM	D	7	3	23,24,25	0.64	0	30,32,34	0.96	2 (6%)							

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	WCM	D	8	3	-	2/20/22/24	0/2/2/2
3	WCM	D	7	3	-	0/20/22/24	0/2/2/2

There are no bond length outliers.

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

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
3	D	8	WCM	C65-C10-C11	-3.98	99.72	109.27
3	D	8	WCM	C13-C12-N6	3.95	116.10	110.64
3	D	8	WCM	O12-C11-C10	-3.01	114.17	119.66



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Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
3	D	8	WCM	C64-N6-C12	2.61	121.28	116.17
3	D	7	WCM	C64-N6-C12	2.46	120.99	116.17

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	D	8	WCM	C10-C65-C66-C71
3	D	8	WCM	C10-C65-C66-C67

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 1 ligands modelled in this entry, 1 is monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	$OWAB(Å^2)$	Q<0.9
1	А	92/124~(74%)	0.07	3 (3%) 46 43	34, 45, 66, 80	0
2	В	471/546 (86%)	0.30	30 (6%) 19 16	28, 47, 85, 112	0
3	D	9/13~(69%)	-0.05	0 100 100	36, 41, 49, 60	0
All	All	572/683~(83%)	0.26	33 (5%) 23 20	28, 47, 84, 112	0

The worst 5 of 33 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	В	570	ASP	5.8
2	В	641	THR	5.0
2	В	640	GLY	4.4
2	В	615	ILE	4.3
2	В	620	GLU	3.9

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

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	$\mathbf{B} ext{-factors}(\mathrm{\AA}^2)$	Q<0.9
3	WCM	D	8	23/24	0.97	0.16	29,38,47,50	0
3	WCM	D	7	23/24	0.98	0.15	29,36,41,42	0

6.3 Carbohydrates (i)

There are no monosaccharides in this entry.



6.4 Ligands (i)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	$\mathbf{B} ext{-factors}(\mathrm{\AA}^2)$	Q<0.9
4	CA	В	701	1/1	0.94	0.06	$65,\!65,\!65,\!65$	0

6.5 Other polymers (i)

There are no such residues in this entry.

