



wwPDB X-ray Structure Validation Summary Report ⓘ

Dec 12, 2016 – 11:03 AM EST

PDB ID : 5JBB
Title : Crystal structure of factor IXa variant V16I K98T Y177T I213V in complex with EGR-chloromethylketone
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Deposited on : 2016-04-13
Resolution : 1.56 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 : unknown
Xtriage (Phenix) : 1.9-1692
EDS : rb-20028442
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) : rb-20028442

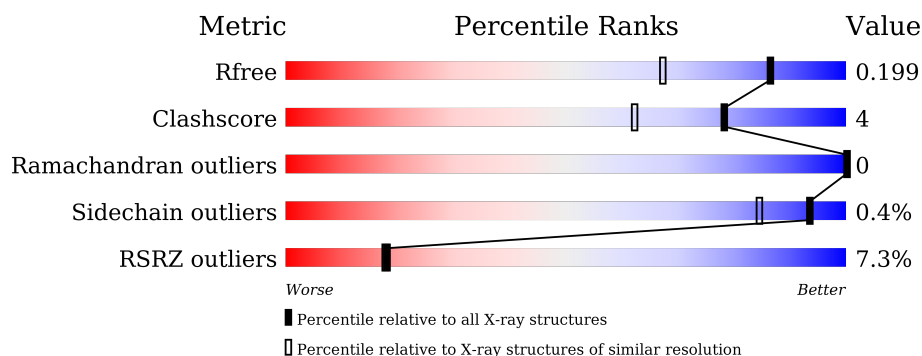
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.56 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	1665 (1.58-1.54)
Clashscore	102246	1014 (1.56-1.56)
Ramachandran outliers	100387	1704 (1.58-1.54)
Sidechain outliers	100360	1702 (1.58-1.54)
RSRZ outliers	91569	1668 (1.58-1.54)

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	E	58	<div> <div>24%</div> <div> <div></div> <div>76%</div> <div>12%</div> <div>12%</div> </div> </div>
2	S	235	<div> <div>3%</div> <div> <div></div> <div>93%</div> <div>7%</div> </div> </div>

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 4771 atoms, of which 2219 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 Coagulation factor IX.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	E	51	Total	C	H	N	O	S	0	0	0
			746	230	363	71	75	7			

- Molecule 2 is a protein called Coagulation factor IX.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	S	235	Total	C	H	N	O	S	0	6	0
			3694	1185	1824	323	349	13			

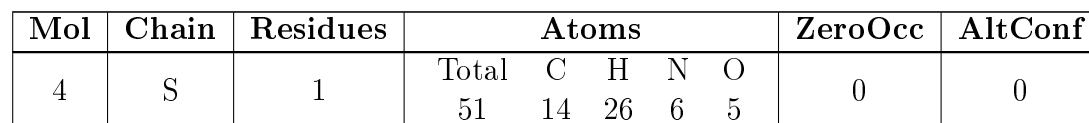
There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
S	16	ILE	VAL	engineered mutation	UNP P00740
S	98	THR	LYS	engineered mutation	UNP P00740
S	177	THR	TYR	engineered mutation	UNP P00740
S	213	VAL	ILE	engineered mutation	UNP P00740

- Molecule 3 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	S	1	Total	Ca	0	0
			1	1		

- Molecule 4 is L-alpha-glutamyl-N-{(1S)-4-[[amino(iminio)methyl]amino]-1-[(1S)-2-chloro-1-hydroxyethyl]butyl}glycinamide (three-letter code: 0GJ) (formula: C₁₄H₂₈ClN₆O₅).



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- The chemical structure of Dimethyl Sulfide (DMS) is shown. It consists of a central sulfur atom (S) bonded to two methyl groups (C1 and C2) and one oxygen atom (O). The sulfur atom is yellow, the carbon atoms are green, and the oxygen atom is red. The bonds are colored: gray for C1-S, yellow for C2-S, and red for S-O. The oxygen atom is labeled with a green 'O'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	S	1	Total	C	H	O	S	0	0
			10	2	6	1	1		

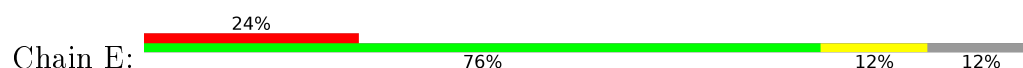
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	E	28	Total 28	O 28	0	0
6	S	241	Total 241	O 241	0	0

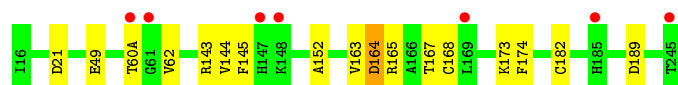
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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: Coagulation factor IX



- Molecule 2: Coagulation factor IX



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	44.31Å 66.61Å 97.05Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	21.60 – 1.56 21.60 – 1.56	Depositor EDS
% Data completeness (in resolution range)	99.7 (21.60-1.56) 99.7 (21.60-1.56)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.73 (at 1.56Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.158 , 0.190 0.169 , 0.199	Depositor DCC
R_{free} test set	2062 reflections (5.22%)	DCC
Wilson B-factor (Å ²)	20.4	Xtriage
Anisotropy	0.176	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 48.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	4771	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.07% 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](#)

Bond lengths and bond angles in the following residue types are not validated in this section: CA, DMS, OGJ

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	E	0.59	0/388	0.67	0/521
2	S	0.80	0/1948	0.83	3/2640 (0.1%)
All	All	0.77	0/2336	0.80	3/3161 (0.1%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	S	21	ASP	CB-CG-OD1	7.44	125.00	118.30
2	S	189	ASP	CB-CG-OD1	5.49	123.25	118.30
2	S	21	ASP	CB-CG-OD2	-5.13	113.69	118.30

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	E	383	363	363	5	0
2	S	1870	1824	1791	11	0
3	S	1	0	0	0	0
4	S	25	26	24	0	0
5	S	4	6	6	2	0
6	E	28	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	S	241	0	0	2	0
All	All	2552	2219	2184	16	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 4.

The worst 5 of 16 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:E:88:CYS:N	1:E:101:ASN:HD21	1.96	0.64
1:E:116:ARG:NH1	1:E:125:GLU:OE1	2.36	0.48
2:S:144:VAL:HG23	2:S:145:PHE:CD2	2.49	0.48
2:S:163:VAL:HB	2:S:182[C]:CYS:SG	2.54	0.48
1:E:116:ARG:HG3	1:E:127:ALA:HB2	1.96	0.47

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	E	49/58 (84%)	45 (92%)	4 (8%)	0	100	100
2	S	243/235 (103%)	236 (97%)	7 (3%)	0	100	100
All	All	292/293 (100%)	281 (96%)	11 (4%)	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	Percentiles	
1	E	45/52 (86%)	45 (100%)	0	100	100
2	S	206/198 (104%)	205 (100%)	1 (0%)	92	81
All	All	251/250 (100%)	250 (100%)	1 (0%)	93	84

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

Mol	Chain	Res	Type
2	S	164	ASP

Some 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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 3 ligands modelled in this entry, 1 is monoatomic - leaving 2 for Mogul analysis.

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).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	0GJ	S	302	2	20,24,25	2.68	4 (20%)	23,30,31	1.62	6 (26%)
5	DMS	S	303	-	3,3,3	0.68	0	3,3,3	1.16	0

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
4	0GJ	S	302	2	-	0/27/29/31	0/0/0/0
5	DMS	S	303	-	-	0/0/0/0	0/0/0/0

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	S	302	0GJ	C-N1	3.19	1.40	1.33
4	S	302	0GJ	C1-N2	5.16	1.44	1.34
4	S	302	0GJ	CZ-NH1	6.22	1.44	1.32
4	S	302	0GJ	CZ-NE	7.24	1.46	1.33

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

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	S	302	0GJ	CD1-NE-CZ	-3.59	116.84	124.68
4	S	302	0GJ	O1-C1-N2	-2.94	118.01	122.96
4	S	302	0GJ	CG1-CD1-NE	-2.89	104.11	112.29
4	S	302	0GJ	CG-CB-CA	-2.29	108.77	114.18
4	S	302	0GJ	CG1-CB1-CA2	-2.14	109.51	114.03

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	S	303	DMS	2	0

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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	E	51/58 (87%)	1.43	14 (27%) 1 1	23, 42, 88, 113	0
2	S	235/235 (100%)	0.00	7 (2%) 54 58	13, 22, 56, 127	0
All	All	286/293 (97%)	0.26	21 (7%) 18 18	13, 25, 70, 127	0

The worst 5 of 21 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	138	SER	7.2
2	S	61	GLY	6.7
1	E	104	ASP	6.5
1	E	103	ALA	5.4
1	E	102	SER	5.0

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](#)

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th 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	LLDF	B-factors(\AA^2)	Q<0.9
5	DMS	S	303	4/4	0.95	0.11	0.91	42,56,60,60	0
4	OGJ	S	302	25/26	0.95	0.09	-0.02	14,20,44,68	0
3	CA	S	301	1/1	1.00	0.04	-2.48	18,18,18,18	0

6.5 Other polymers [i](#)

There are no such residues in this entry.