

Full wwPDB X-ray Structure Validation Report (i)

Feb 25, 2024 – 03:17 AM EST

PDB ID : 5E67

Title: K103A/K262A double mutant of I-SmaMI

Authors: Shen, B.W.; Stoddard, B.

Deposited on : 2015-10-09

Resolution : 2.20 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
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 : 2.36

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)

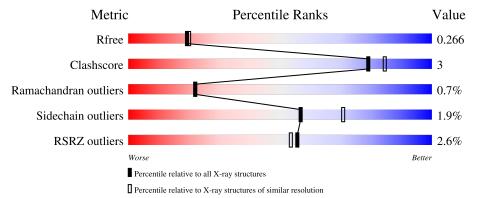
Validation Pipeline (wwPDB-VP) : 2.36

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 2.20 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 $(\# \mathrm{Entries})$	$\begin{array}{c} {\rm Similar\ resolution} \\ (\#{\rm Entries},\ {\rm resolution\ range}({\rm \AA})) \end{array}$		
R_{free}	130704	4898 (2.20-2.20)		
Clashscore	141614	5594 (2.20-2.20)		
Ramachandran outliers	138981	5503 (2.20-2.20)		
Sidechain outliers	138945	5504 (2.20-2.20)		
RSRZ outliers	127900	4800 (2.20-2.20)		

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	302	86%		10% • •			
2	В	25	64%	36%				
3	С	25	56%	32%	12%			



2 Entry composition (i)

There are 7 unique types of molecules in this entry. The entry contains 3578 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 I-SmaMI LAGLIDADG meganuclease.

\mathbf{Mol}	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	A	294	Total 2348	C 1532	N 386	O 422	S 8	0	1	0

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	103	ALA	LYS	engineered mutation	UNP F7WD42
A	165	ASN	LEU	engineered mutation	UNP F7WD42
A	262	ALA	LYS	engineered mutation	UNP F7WD42
A	267	GLN	MET	engineered mutation	UNP F7WD42

• Molecule 2 is a DNA chain called DNA bottom strand.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
2	R	25	Total	С	N	О	Р	0	0	0
	D	20	511	244	98	145	24	0		

• Molecule 3 is a DNA chain called DNA top strand.

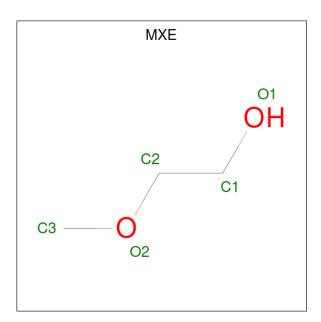
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
2	C	25	Total	С	N	О	Р	0	0	0
)		2.9	508	244	89	151	24	U		

• Molecule 4 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	2	Total Mg 2 2	0	0

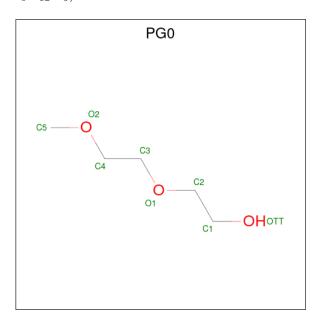
• Molecule 5 is 2-METHOXYETHANOL (three-letter code: MXE) (formula: $C_3H_8O_2$).





Mol	Chain	Residues	Aton	ns	ZeroOcc	AltConf
5	A	1	Total (C O 3 2	0	0

 \bullet Molecule 6 is 2-(2-METHOXYETHOXY) ETHANOL (three-letter code: PG0) (formula: $C_5H_{12}O_3).$



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	В	1	Total 16	C 10	O 6	0	1

• Molecule 7 is water.



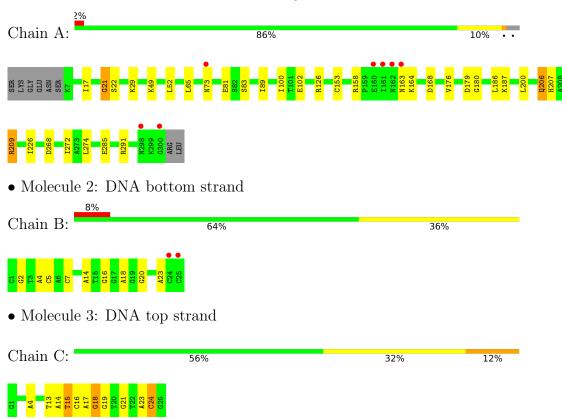
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	133	Total O 133 133	0	0
7	В	27	Total O 27 27	0	0
7	С	28	Total O 28 28	0	0



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: I-SmaMI LAGLIDADG meganuclease





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants	60.52Å 67.33Å 98.57Å	Depositor
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor
Resolution (Å)	25.00 - 2.20	Depositor
rtesolution (A)	24.08 - 2.20	EDS
% Data completeness	92.7 (25.00-2.20)	Depositor
(in resolution range)	92.8 (24.08-2.20)	EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$< I/\sigma(I) > 1$	2.11 (at 2.19Å)	Xtriage
Refinement program	REFMAC 5.8.0124	Depositor
P. P.	0.179 , 0.254	Depositor
R, R_{free}	0.188 , 0.266	DCC
R_{free} test set	957 reflections (4.90%)	wwPDB-VP
Wilson B-factor (Å ²)	33.2	Xtriage
Anisotropy	0.096	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.33, 36.4	EDS
L-test for twinning ²	$ < L >=0.50, < L^2>=0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	3578	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 5.92% 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: MG, MXE, PG0

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).

Mal	Chain	Bo	nd lengths	Bond angles		
Mol		RMSZ	# Z > 5	RMSZ	# Z > 5	
1	A	1.08	4/2402 (0.2%)	1.06	8/3234 (0.2%)	
2	В	0.87	1/574~(0.2%)	1.29	7/884 (0.8%)	
3	С	0.74	1/568~(0.2%)	1.14	5/875~(0.6%)	
All	All	1.00	$6/3544 \ (0.2\%)$	1.12	$20/4993 \ (0.4\%)$	

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$\operatorname{Observed}(\text{\AA})$	$\operatorname{Ideal}(\text{\AA})$
2	В	18	DA	O3'-P	-6.76	1.53	1.61
1	A	21	GLY	N-CA	6.55	1.55	1.46
1	A	285	GLU	CD-OE1	5.70	1.31	1.25
1	A	83	SER	CB-OG	-5.43	1.35	1.42
1	A	285	GLU	CG-CD	5.29	1.59	1.51
3	С	17	DA	O3'-P	-5.05	1.55	1.61

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^o)$	$\operatorname{Ideal}({}^{o})$
2	В	5	DC	O5'-P-OP2	-13.96	93.13	105.70
2	В	5	DC	O5'-P-OP1	10.19	122.93	110.70
3	С	15	DT	O5'-P-OP1	-7.53	98.92	105.70
2	В	16	DG	C1'-O4'-C4'	-7.34	102.76	110.10
2	В	5	DC	C1'-O4'-C4'	-7.21	102.89	110.10
3	С	24	DC	O5'-P-OP2	-6.70	99.67	105.70
3	С	18	DG	C1'-O4'-C4'	-6.61	103.50	110.10
3	С	16	DC	O5'-P-OP1	-6.45	99.90	105.70
1	A	126	ARG	NE-CZ-NH1	6.01	123.31	120.30
1	A	168	ASP	CB-CG-OD1	6.01	123.71	118.30
1	A	52	LEU	CA-CB-CG	6.00	129.09	115.30
1	A	268	ASP	CB-CG-OD2	5.97	123.67	118.30

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Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$\mathbf{Observed}(^o)$	$\operatorname{Ideal}({}^{o})$
1	A	81	GLU	OE1-CD-OE2	-5.89	116.23	123.30
2	В	7	DC	C1'-O4'-C4'	-5.75	104.35	110.10
2	В	20	DG	O5'-P-OP2	-5.68	100.59	105.70
1	A	226	ILE	CG1-CB-CG2	-5.59	99.11	111.40
2	В	2	DG	C1'-O4'-C4'	-5.45	104.66	110.10
3	С	13	DT	C4'-C3'-O3'	5.42	123.26	109.70
1	A	158	ARG	NE-CZ-NH2	-5.22	117.69	120.30
1	A	209	ARG	NE-CZ-NH2	-5.18	117.71	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	A	2348	0	2405	11	0
2	В	511	0	282	3	0
3	С	508	0	285	5	0
4	A	2	0	0	0	0
5	A	5	0	8	1	0
6	В	16	0	24	1	0
7	A	133	0	0	0	0
7	В	27	0	0	1	0
7	С	28	0	0	0	0
All	All	3578	0	3004	18	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 3.

All (18) 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)
3:C:14:DA:H1'	3:C:15:DT:O4'	2.03	0.58
1:A:207:HIS:HD2	1:A:209:ARG:H	1.56	0.54
2:B:4:DA:OP1	7:B:201:HOH:O	2.19	0.53

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Atom-1	Atom-2	Interatomic	Clash
Atom-1	Atom-2	${ m distance}({ m \AA})$	overlap (Å)
3:C:18:DG:H2"	3:C:19:DG:O5'	2.09	0.52
5:A:403:MXE:H12	2:B:14:DA:OP1	2.10	0.51
1:A:100:ILE:O	1:A:164[A]:LYS:HE3	2.12	0.50
2:B:23:DA:C2	3:C:4:DA:C2	3.01	0.49
1:A:17:ILE:HG23	1:A:21:GLY:HA3	1.95	0.48
1:A:176:VAL:HG13	1:A:180:GLY:HA3	1.97	0.47
1:A:187:LYS:HE3	3:C:21:DG:O6	2.16	0.46
1:A:102:GLU:HG3	1:A:209:ARG:CD	2.47	0.44
3:C:23:DA:H2"	3:C:24:DC:O5'	2.17	0.44
1:A:65:LEU:HD12	1:A:89:ILE:HG13	2.00	0.44
1:A:179:ASP:OD2	1:A:206:GLN:OE1	2.35	0.43
6:B:101[A]:PG0:H31	6:B:101[A]:PG0:H53	1.74	0.41
1:A:274:LEU:HD12	1:A:274:LEU:HA	1.91	0.41
1:A:200:LEU:HD21	1:A:272:ILE:CG2	2.50	0.41
1:A:186:LEU:CD1	1:A:291:ARG:HG2	2.52	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	A	293/302 (97%)	284 (97%)	7 (2%)	2 (1%)	22 22

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	153	CYS
1	A	163	ASN



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.

\mathbf{N}	ſol	Chain	Analysed	Rotameric	Outliers	Percentiles
	1	A	258/264 (98%)	253 (98%)	5 (2%)	57 71

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

Mol	Chain	Res	Type
1	A	22	SER
1	A	29	LYS
1	A	49	LYS
1	A	73	ASN
1	A	206	GLN

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

Mol	Chain	Res	Type
1	A	85	GLN
1	A	207	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)

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)

Of 5 ligands modelled in this entry, 2 are monoatomic - leaving 3 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		pe Chain	hain Res	Res Link	Bond lengths			Bond angles		
MIOI	Type	Chain	nes	LIIIK	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	PG0	В	101[B]	-	7,7,7	0.53	0	6,6,6	0.67	0
6	PG0	В	101[A]	-	7,7,7	0.63	0	6,6,6	0.57	0
5	MXE	A	403	-	4,4,4	0.74	0	3,3,3	0.68	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
6	PG0	В	101[B]	-	-	5/5/5/5	-
6	PG0	В	101[A]	-	-	4/5/5/5	-
5	MXE	A	403	-	-	2/2/2/2	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	В	101[A]	PG0	OTT-C1-C2-O1
6	В	101[A]	PG0	O1-C3-C4-O2
6	В	101[B]	PG0	O1-C3-C4-O2
5	A	403	MXE	O1-C1-C2-O2
5	A	403	MXE	C1-C2-O2-C3
6	В	101[A]	PG0	C3-C4-O2-C5
6	В	101[B]	PG0	C1-C2-O1-C3
6	В	101[A]	PG0	C1-C2-O1-C3
6	В	101[B]	PG0	OTT-C1-C2-O1

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Mol	Chain	Res	Type	Atoms
6	В	101[B]	PG0	C4-C3-O1-C2
6	В	101[B]	PG0	C3-C4-O2-C5

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	В	101[A]	PG0	1	0
5	A	403	MXE	1	0

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	$\mathbf{OWAB}(\mathrm{\AA}^2)$	Q < 0.9
1	A	294/302~(97%)	-0.20	7 (2%) 59 56	16, 29, 52, 79	2 (0%)
2	В	25/25 (100%)	-0.20	2 (8%) 12 11	25, 39, 59, 94	0
3	С	25/25 (100%)	-0.58	0 100 100	32, 42, 52, 55	0
All	All	344/352 (97%)	-0.23	9 (2%) 56 53	16, 31, 54, 94	2 (0%)

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	163	ASN	5.3
2	В	25	DC	2.9
1	A	162	ASN	2.8
1	A	300	GLY	2.5
1	A	161	ILE	2.5
1	A	298	ASN	2.4
2	В	24	DC	2.4
1	A	160	GLU	2.3
1	A	73	ASN	2.2

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)

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	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{\mathring{A}}^2)$	Q < 0.9
5	MXE	A	403	5/5	0.73	0.23	47,50,52,55	0
6	PG0	В	101[A]	8/8	0.77	0.27	21,30,36,37	8
6	PG0	В	101[B]	8/8	0.77	0.27	30,40,49,50	8
4	MG	A	402	1/1	0.89	0.10	48,48,48,48	0
4	MG	A	401	1/1	0.94	0.08	41,41,41,41	0

6.5 Other polymers (i)

There are no such residues in this entry.

