

wwPDB X-ray Structure Validation Summary Report (i)

Nov 2, 2021 – 07:10 AM EDT

PDB ID : 3L6C

Title: X-ray crystal structure of rat serine racemase in complex with malonate a

potent inhibitor

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Deposited on : 2009-12-23

Resolution : 2.20 Å(reported)

This is a wwPDB X-ray Structure Validation Summary Report for a publicly released PDB entry.

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 : 2.23.2

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)

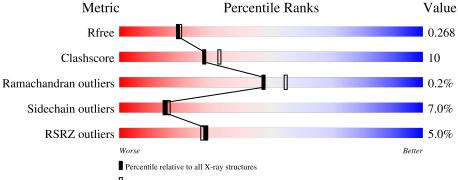
Validation Pipeline (wwPDB-VP) : 2.23.2

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 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.



 $\ensuremath{\overline{\square}}$ Percentile relative to X-ray structures of similar resolution

Metric	$\begin{array}{c} \text{Whole archive} \\ (\#\text{Entries}) \end{array}$	Similar resolution $(\#\text{Entries, resolution range}(\mathring{A}))$		
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	339	80%		13%	• 5%
1	В	339	65%	24%	•	8%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:



Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	MLI	В	341	_	_	X	-



2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 4905 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 Serine racemase.

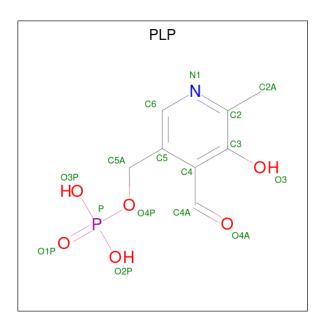
\mathbf{Mol}	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	Δ	322	Total	С	N	О	S	0	0	0
	Λ	322	2423	1541	407	469	6		U	
1	B	312	Total	С	N	О	S	0	0	0
1	Ъ	312	2356	1501	397	452	6			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	2	ASP	CYS	engineered mutation	UNP Q76EQ0
A	6	ASP	CYS	engineered mutation	UNP Q76EQ0
A	334	HIS	-	expression tag	UNP Q76EQ0
A	335	HIS	-	expression tag	UNP Q76EQ0
A	336	HIS	-	expression tag	UNP Q76EQ0
A	337	HIS	-	expression tag	UNP Q76EQ0
A	338	HIS	-	expression tag	UNP Q76EQ0
A	339	HIS	-	expression tag	UNP Q76EQ0
В	2	ASP	CYS	engineered mutation	UNP Q76EQ0
В	6	ASP	CYS	engineered mutation	UNP Q76EQ0
В	334	HIS	-	expression tag	UNP Q76EQ0
В	335	HIS	-	expression tag	UNP Q76EQ0
В	336	HIS	-	expression tag	UNP Q76EQ0
В	337	HIS	-	expression tag	UNP Q76EQ0
В	338	HIS	-	expression tag	UNP Q76EQ0
В	339	HIS	-	expression tag	UNP Q76EQ0

• Molecule 2 is PYRIDOXAL-5'-PHOSPHATE (three-letter code: PLP) (formula: C₈H₁₀NO₆P).





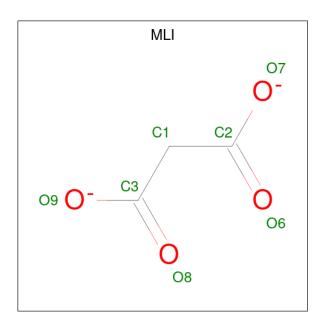
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf		
9	٨	1	Total	С	N	О	Р	0	0	
2	A	1	15	8	1	5	1	0	0	
9	D	1	Total	С	N	О	Р	0	0	
2	Б	1	15	8	1	5	1	0	U	

• Molecule 3 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

I	Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
	3	A	1	Total Mn 1 1	0	0
	3	В	1	Total Mn 1 1	0	0

 \bullet Molecule 4 is MALONATE ION (three-letter code: MLI) (formula: $\mathrm{C_3H_2O_4}).$





Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C O 7 3 4	0	0
4	В	1	Total C O 7 3 4	0	0

• Molecule 5 is water.

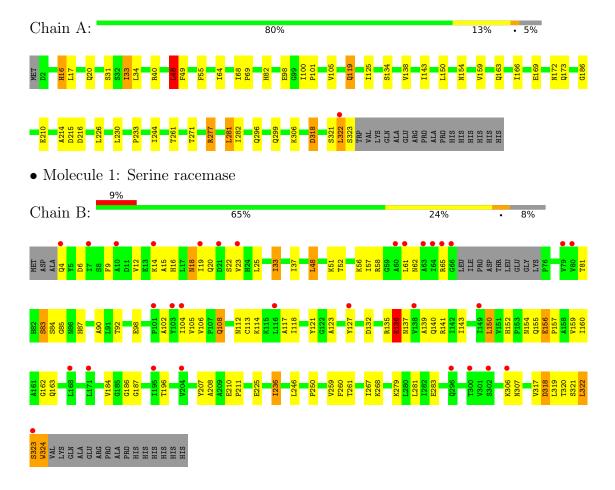
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	62	Total O 62 62	0	0
5	В	18	Total O 18 18	0	0



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: Serine racemase





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 3	Depositor
Cell constants	110.59Å 110.59Å 47.92Å	Donositon
a, b, c, α , β , γ	90.00° 90.00° 120.00°	Depositor
Resolution (Å)	47.89 - 2.20	Depositor
rtesolution (A)	28.88 - 2.20	EDS
% Data completeness	95.9 (47.89-2.20)	Depositor
(in resolution range)	95.9 (28.88-2.20)	EDS
R_{merge}	0.06	Depositor
R_{sym}	0.06	Depositor
$< I/\sigma(I) > 1$	1.76 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.216 , 0.267	Depositor
it, it free	0.221 , 0.268	DCC
R_{free} test set	1586 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	39.5	Xtriage
Anisotropy	0.080	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.37, 33.7	EDS
L-test for twinning ²	$< L > = 0.48, < L^2> = 0.31$	Xtriage
	0.033 for -h,-k,l	
Estimated twinning fraction	0.079 for h,-h-k,-l	Xtriage
	0.028 for -k,-h,-l	
F_o, F_c correlation	0.94	EDS
Total number of atoms	4905	wwPDB-VP
Average B, all atoms (\mathring{A}^2)	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 4.07% 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: PLP, MN, MLI

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		
IVIOI		RMSZ	# Z > 5	RMSZ	# Z > 5	
1	A	0.76	$1/2467 \ (0.0\%)$	0.81	3/3365 (0.1%)	
1	В	0.65	0/2400	0.73	$2/3272 \ (0.1\%)$	
All	All	0.71	1/4867 (0.0%)	0.77	5/6637 (0.1%)	

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\operatorname{Observed}(\text{\AA})$	$\operatorname{Ideal}(ext{\AA})$
1	A	119	GLN	CD-OE1	5.52	1.36	1.24

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\mathbf{Observed}(^o)$	$\operatorname{Ideal}(^{o})$
1	В	48	LEU	CA-CB-CG	7.22	131.90	115.30
1	A	216	ASP	CB-CG-OD2	-6.69	112.28	118.30
1	A	48	LEU	CA-CB-CG	5.31	127.51	115.30
1	В	323	SER	N-CA-CB	5.29	118.44	110.50
1	A	216	ASP	CB-CG-OD1	5.17	122.96	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	A	2423	0	2467	35	0
1	В	2356	0	2394	64	0
2	A	15	0	6	0	0
2	В	15	0	6	0	0
3	A	1	0	0	0	0
3	В	1	0	0	0	0
4	A	7	0	2	0	0
4	В	7	0	2	5	0
5	A	62	0	0	0	0
5	В	18	0	0	0	0
All	All	4905	0	4877	96	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 10.

The worst 5 of 96 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$\begin{array}{c} {\rm Interatomic} \\ {\rm distance} \ ({\rm \AA}) \end{array}$	$\begin{array}{c} \text{Clash} \\ \text{overlap } (\text{\AA}) \end{array}$
1:B:154:ASN:ND2	1:B:186:GLY:HA3	1.12	1.41
1:B:154:ASN:ND2	1:B:186:GLY:CA	2.06	1.18
1:B:154:ASN:HD22	1:B:186:GLY:CA	1.63	1.09
1:A:322:LEU:C	1:A:322:LEU:HD23	1.70	1.09
1:A:322:LEU:HG	1:A:323:SER:N	1.68	1.06

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	Perce	entiles
1	A	320/339 (94%)	304 (95%)	16 (5%)	0	100	100
1	В	308/339 (91%)	279 (91%)	28 (9%)	1 (0%)	41	46
All	All	628/678 (93%)	583 (93%)	44 (7%)	1 (0%)	47	55



All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	В	136	GLU

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	A	268/283 (95%)	258 (96%)	10 (4%)	34 43		
1	В	260/283 (92%)	233 (90%)	27 (10%)	7 6		
All	All	528/566 (93%)	491 (93%)	37 (7%)	15 16		

5 of 37 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	В	157	PRO
1	В	322	LEU
1	В	163	GLN
1	В	306	LYS
1	В	9	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	36	GLN
1	A	137	ASN
1	A	229	ASN
1	В	307	ASN

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 6 ligands modelled in this entry, 2 are monoatomic - leaving 4 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	Trimo	Chain	Res	es Link	Bond lengths			Bond angles		
MIOI	Type	Chain	nes	Lilik	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	MLI	A	341	-	0,6,6	-	-	0,7,7	-	-
2	PLP	A	350	1	15,15,16	1.13	1 (6%)	20,22,23	1.75	6 (30%)
2	PLP	В	350	1	15,15,16	0.88	1 (6%)	20,22,23	1.55	4 (20%)
4	MLI	В	341	-	0,6,6	-	-	0,7,7	-	-

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	MLI	A	341	-	-	0/0/4/4	-
2	PLP	A	350	1	-	0/6/6/8	0/1/1/1
2	PLP	В	350	1	-	3/6/6/8	0/1/1/1
4	MLI	В	341	-	-	0/0/4/4	-

All (2) bond length outliers are listed below:

	Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$\operatorname{Observed}(\mathring{\mathbf{A}})$	$\operatorname{Ideal}(\text{\AA})$
	2	A	350	PLP	C3-C2	-2.67	1.38	1.40
Ī	2	В	350	PLP	C3-C2	-2.54	1.38	1.40



The worst 5 of 10 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^o)$	$Ideal(^{o})$
2	В	350	PLP	O4P-C5A-C5	3.27	115.58	109.35
2	A	350	PLP	C4A-C4-C5	3.19	124.23	120.94
2	A	350	PLP	C2A-C2-N1	3.12	123.77	117.67
2	A	350	PLP	O4P-C5A-C5	3.12	115.29	109.35
2	A	350	PLP	C4A-C4-C3	-2.85	115.66	120.50

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	В	350	PLP	C4-C5-C5A-O4P
2	В	350	PLP	C6-C5-C5A-O4P
2	В	350	PLP	C5A-O4P-P-O1P

There are no ring outliers.

1 monomer is involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	В	341	MLI	5	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	$\langle { m RSRZ} \rangle$	$\# \mathrm{RSRZ}{>}2$	$OWAB(A^2)$	Q<0.9
1	A	322/339~(94%)	-0.33	1 (0%) 94 93	20, 33, 50, 64	0
1	В	312/339 (92%)	0.48	31 (9%) 7 6	31, 62, 82, 85	0
All	All	634/678 (93%)	0.07	32 (5%) 28 27	20, 43, 80, 85	0

The worst 5 of 32 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	322	LEU	5.1
1	В	19	ILE	4.8
1	В	66	GLY	4.3
1	В	138	VAL	4.0
1	В	7	ILE	3.8

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)

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	\mathbf{Type}	Chain	Res	Atoms	RSCC	RSR	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{\mathring{A}}^2)$	Q<0.9
3	MN	В	340	1/1	0.71	0.08	60,60,60,60	1
4	MLI	В	341	7/7	0.88	0.12	44,47,50,51	0
3	MN	A	340	1/1	0.90	0.06	45,45,45,45	1
4	MLI	A	341	7/7	0.97	0.11	25,30,35,36	0
2	PLP	В	350	15/16	0.97	0.10	38,40,42,44	0
2	PLP	A	350	15/16	0.99	0.11	20,27,30,30	0

6.5 Other polymers (i)

There are no such residues in this entry.

