



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 26, 2014 – 03:29 PM GMT

PDB ID : 3L6C  
Title : X-ray crystal structure of rat serine racemase in complex with malonate a potent inhibitor  
Authors : Smith, M.A.; Mack, V.; Ebner, A.; Moraes, I.; Felicetti, B.; Wood, M.; Schonfeld, D.; Mather, O.; Cesura, A.; Barker, J.  
Deposited on : 2009-12-23  
Resolution : 2.20 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.  
We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)  
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

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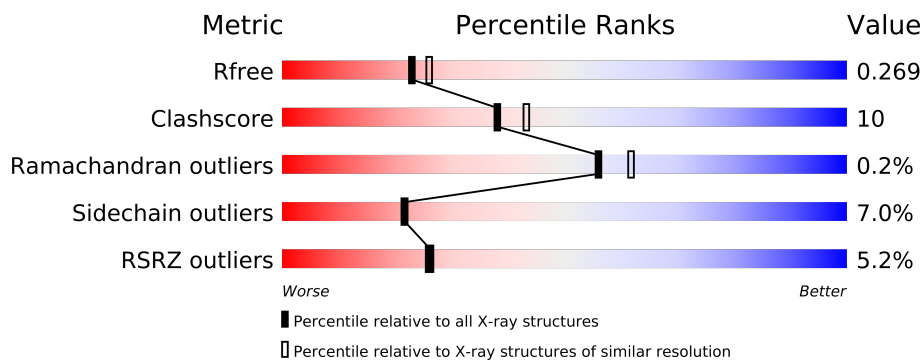
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : dev-1323  
EDS : stable22639  
Percentile statistics : 21963  
Refmac : 5.8.0049  
CCP4 : 6.3.0 (Settle)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance



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 (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	66092	2938 (2.20-2.20)
Clashscore	79885	3751 (2.20-2.20)
Ramachandran outliers	78287	3681 (2.20-2.20)
Sidechain outliers	78261	3682 (2.20-2.20)
RSRZ outliers	66119	2939 (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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	339	
1	B	339	

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 4905 atoms, of which 0 are hydrogen and 0 are deuterium.

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.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	322	Total	C	N	O	S	0	0	0
			2423	1541	407	469	6			
1	B	312	Total	C	N	O	S	0	0	0
			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	UNP Q76EQ0
A	6	ASP	CYS	ENGINEERED	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
B	2	ASP	CYS	ENGINEERED	UNP Q76EQ0
B	6	ASP	CYS	ENGINEERED	UNP Q76EQ0
B	334	HIS	-	EXPRESSION TAG	UNP Q76EQ0
B	335	HIS	-	EXPRESSION TAG	UNP Q76EQ0
B	336	HIS	-	EXPRESSION TAG	UNP Q76EQ0
B	337	HIS	-	EXPRESSION TAG	UNP Q76EQ0
B	338	HIS	-	EXPRESSION TAG	UNP Q76EQ0
B	339	HIS	-	EXPRESSION TAG	UNP Q76EQ0

- Molecule 2 is PYRIDOXAL-5'-PHOSPHATE (three-letter code: PLP) (formula: C<sub>8</sub>H<sub>10</sub>NO<sub>6</sub>P).

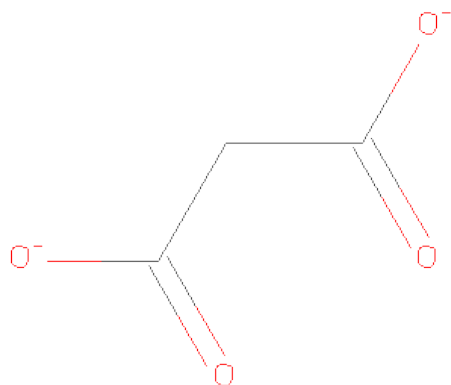


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			15	8	1	5	1		
2	B	1	Total	C	N	O	P	0	0
			15	8	1	5	1		

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

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

- Molecule 4 is MALONATE ION (three-letter code: MLI) (formula: C<sub>3</sub>H<sub>2</sub>O<sub>4</sub>).



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

- Molecule 5 is water.

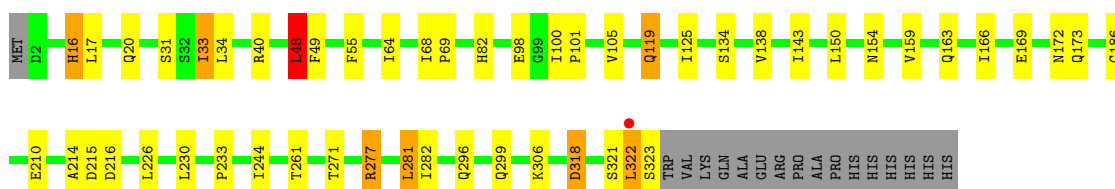
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	62	Total	O	0	0
			62	62		
5	B	18	Total	O	0	0
			18	18		

### 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 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.

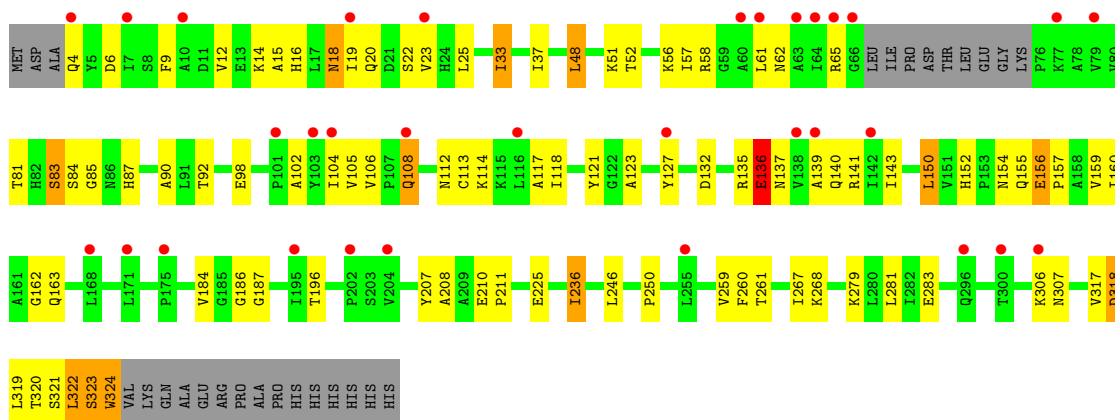
- Molecule 1: Serine racemase

Chain A:



- Molecule 1: Serine racemase

Chain B:



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 3	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	110.59Å 110.59Å 47.92Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	47.89 – 2.20 28.88 – 2.20	Depositor EDS
% Data completeness (in resolution range)	95.9 (47.89-2.20) 95.9 (28.88-2.20)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	0.06	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.76 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, $R_{free}$	0.216 , 0.267 0.222 , 0.269	Depositor DCC
$R_{free}$ test set	1586 reflections (5.24%)	DCC
Wilson B-factor (Å <sup>2</sup> )	39.5	Xtriage
Anisotropy	0.080	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.37 , 36.3	EDS
Estimated twinning fraction	0.033 for -h,-k,l 0.079 for h,-h-k,-l 0.028 for -k,-h,-l	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 31868 reflections	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	4905	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	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.*

<sup>1</sup>Intensities estimated from amplitudes.

## 5 Model quality ⓘ

### 5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: MLI, MN, PLP

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	A	0.76	1/2467 (0.0%)	0.81	3/3365 (0.1%)
1	B	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	Observed(Å)	Ideal(Å)
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	Observed(°)	Ideal(°)
1	B	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	B	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 Close contacts ⓘ

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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.



Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2423	0	2467	35	0
1	B	2356	0	2394	64	0
2	A	15	0	6	0	0
2	B	15	0	6	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	7	0	2	0	0
4	B	7	0	2	5	0
5	A	62	0	0	0	0
5	B	18	0	0	0	0
All	All	4905	0	4877	96	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 10.

All (96) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
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:HD23	1:A:322:LEU:C	1.70	1.09
1:A:322:LEU:HG	1:A:323:SER:N	1.68	1.06
1:B:236:ILE:HD13	1:B:268:LYS:HE2	1.42	1.01
1:A:322:LEU:HG	1:A:323:SER:H	1.34	0.92
1:B:154:ASN:HD21	1:B:186:GLY:HA3	1.23	0.91
1:A:322:LEU:CG	1:A:323:SER:N	2.32	0.89
1:B:143:ILE:HD11	1:B:150:LEU:HB2	1.56	0.85
1:B:281:LEU:HD21	1:B:319:LEU:HD21	1.61	0.83
1:A:322:LEU:CD2	1:A:322:LEU:C	2.41	0.82
1:B:322:LEU:HB2	1:B:324:TRP:CD1	2.14	0.82
1:B:117:ALA:O	1:B:121:TYR:HD2	1.65	0.78
1:B:15:ALA:O	1:B:19:ILE:HG12	1.85	0.76
1:A:322:LEU:HD23	1:A:322:LEU:O	1.88	0.74
1:B:132:ASP:O	1:B:136:GLU:HB2	1.89	0.73
1:B:236:ILE:HD11	1:B:324:TRP:NE1	2.09	0.68
1:B:108:GLN:HG2	1:B:127:TYR:HB3	1.76	0.68
1:B:322:LEU:CB	1:B:324:TRP:CD1	2.77	0.67
1:B:14:LYS:O	1:B:18:ASN:HB2	1.97	0.64
1:B:104:ILE:HG22	1:B:106:VAL:HG23	1.78	0.64
1:B:154:ASN:HD22	1:B:186:GLY:HA3	0.80	0.63
1:A:134:SER:O	1:A:138:VAL:HG23	2.01	0.61
1:B:207:TYR:CD1	1:B:260:PHE:HE2	2.18	0.61
1:A:226:LEU:CD1	1:A:244:ILE:HG13	2.32	0.59

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:210:GLU:O	1:B:261:THR:HA	2.02	0.58
1:B:281:LEU:HD21	1:B:319:LEU:CD2	2.34	0.57
1:B:211:PRO:HB3	1:B:267:ILE:HD11	1.85	0.57
1:B:83:SER:OG	4:B:341:MLI:O9	2.17	0.57
1:A:318:ASP:HB2	1:A:321:SER:HB3	1.87	0.56
1:B:184:VAL:HG23	1:B:187:GLY:HA2	1.87	0.56
1:B:318:ASP:C	1:B:320:THR:H	2.07	0.56
1:B:321:SER:C	1:B:323:SER:H	2.08	0.55
1:B:104:ILE:HG21	1:B:118:ILE:HD13	1.88	0.55
1:B:52:THR:HG21	1:B:90:ALA:HA	1.87	0.55
1:B:57:ILE:HA	1:B:90:ALA:HB1	1.88	0.55
1:A:33:ILE:HD12	1:A:277:ARG:HG3	1.88	0.55
1:B:51:LYS:O	1:B:52:THR:OG1	2.25	0.54
1:A:119:GLN:HG3	1:A:125:ILE:HD12	1.90	0.54
1:A:322:LEU:CD2	1:A:323:SER:N	2.68	0.54
1:B:56:LYS:CE	4:B:341:MLI:O8	2.56	0.53
1:B:56:LYS:HE3	4:B:341:MLI:O8	2.08	0.53
1:A:296:GLN:O	1:A:299:GLN:HB2	2.09	0.51
1:A:16:HIS:O	1:A:20:GLN:HB3	2.10	0.51
1:B:84:SER:OG	4:B:341:MLI:H12	2.11	0.51
1:B:236:ILE:CD1	1:B:268:LYS:HE2	2.28	0.51
1:B:19:ILE:O	1:B:23:VAL:HB	2.12	0.50
1:A:143:ILE:HD11	1:A:150:LEU:HB2	1.93	0.50
1:A:281:LEU:HD23	1:B:281:LEU:HD13	1.93	0.49
1:B:154:ASN:O	1:B:160:ILE:HG12	2.12	0.49
1:B:321:SER:C	1:B:323:SER:N	2.65	0.49
1:B:65:ARG:NH2	1:B:98:GLU:OE2	2.45	0.49
1:B:318:ASP:O	1:B:320:THR:N	2.44	0.49
1:A:64:ILE:CG2	1:A:98:GLU:HG3	2.43	0.48
1:B:85:GLY:HA3	1:B:114:LYS:HD2	1.94	0.48
1:A:119:GLN:HG3	1:A:125:ILE:CD1	2.44	0.47
1:B:61:LEU:HD22	1:B:65:ARG:HH12	1.80	0.47
1:B:156:GLU:HG2	1:B:159:VAL:HG23	1.97	0.47
1:A:210:GLU:O	1:A:261:THR:HA	2.15	0.47
1:B:208:ALA:HB3	1:B:259:VAL:HG22	1.95	0.46
1:B:16:HIS:O	1:B:20:GLN:N	2.48	0.46
1:A:48:LEU:HD22	1:B:279:LYS:HB2	1.96	0.46
1:A:48:LEU:HD22	1:B:279:LYS:CB	2.46	0.46
1:A:154:ASN:ND2	1:A:186:GLY:HA3	2.31	0.46
1:A:31:SER:HB3	1:A:34:LEU:HB2	1.98	0.46
1:A:271:THR:HA	1:A:282:ILE:HD11	1.98	0.45
1:B:152:HIS:ND1	1:B:155:GLN:HB2	2.32	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:322:LEU:CB	1:B:324:TRP:HD1	2.30	0.44
1:A:169:GLU:O	1:A:173:GLN:HG3	2.16	0.44
1:A:214:ALA:HB2	1:A:233:PRO:HB2	2.00	0.43
1:A:159:VAL:O	1:A:163:GLN:HG2	2.19	0.43
1:B:135:ARG:C	1:B:137:ASN:H	2.21	0.43
1:A:82:HIS:HA	1:A:105:VAL:O	2.18	0.43
1:A:100:ILE:HA	1:A:101:PRO:HD3	1.91	0.43
1:B:92:THR:HG23	1:B:123:ALA:HB2	2.01	0.42
1:B:105:VAL:HG21	1:B:139:ALA:HA	2.00	0.42
1:B:56:LYS:HE2	4:B:341:MLI:O8	2.20	0.42
1:B:12:VAL:HG11	1:B:196:THR:OG1	2.20	0.42
1:A:277:ARG:HA	1:A:277:ARG:HD2	1.80	0.41
1:B:16:HIS:O	1:B:20:GLN:HB3	2.20	0.41
1:A:48:LEU:HD13	1:A:49:PHE:CD1	2.55	0.41
1:A:40:ARG:HA	1:A:306:LYS:HE2	2.02	0.41
1:A:55:PHE:HB2	1:A:166:ILE:HD11	2.02	0.41
1:B:81:THR:HG21	1:B:87:HIS:CD2	2.56	0.41
1:B:246:LEU:O	1:B:250:PRO:HD3	2.21	0.41
1:B:33:ILE:O	1:B:37:ILE:HG13	2.20	0.41
1:B:102:ALA:O	1:B:123:ALA:HB1	2.20	0.41
1:B:322:LEU:HB3	1:B:324:TRP:CD1	2.53	0.41
1:A:210:GLU:CD	1:A:215:ASP:HA	2.41	0.41
1:B:283:GLU:HA	1:B:317:VAL:HB	2.02	0.41
1:B:318:ASP:C	1:B:320:THR:N	2.74	0.41
1:B:58:ARG:HB3	1:B:162:GLY:O	2.21	0.41
1:B:236:ILE:HD11	1:B:324:TRP:HE1	1.85	0.40
1:B:85:GLY:HA3	1:B:114:LYS:CD	2.51	0.40
1:A:68:ILE:HA	1:A:69:PRO:HD3	1.87	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles

### 5.3.1 Protein backbone ⓘ

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	320/339 (94%)	304 (95%)	16 (5%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	308/339 (91%)	279 (91%)	28 (9%)	1 (0%)	50	53
All	All	628/678 (93%)	583 (93%)	44 (7%)	1 (0%)	56	62

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	136	GLU

### 5.3.2 Protein sidechains ⓘ

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%)	45	54
1	B	260/283 (92%)	233 (90%)	27 (10%)	10	9
All	All	528/566 (93%)	491 (93%)	37 (7%)	21	22

All (37) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	16	HIS
1	A	17	LEU
1	A	33	ILE
1	A	48	LEU
1	A	172	ASN
1	A	230	LEU
1	A	277	ARG
1	A	281	LEU
1	A	318	ASP
1	A	322	LEU
1	B	4	GLN
1	B	6	ASP
1	B	9	PHE
1	B	18	ASN
1	B	22	SER
1	B	25	LEU
1	B	33	ILE

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Mol	Chain	Res	Type
1	B	48	LEU
1	B	62	ASN
1	B	83	SER
1	B	108	GLN
1	B	112	ASN
1	B	113	CYS
1	B	136	GLU
1	B	140	GLN
1	B	141	ARG
1	B	150	LEU
1	B	156	GLU
1	B	157	PRO
1	B	163	GLN
1	B	225	GLU
1	B	236	ILE
1	B	306	LYS
1	B	307	ASN
1	B	318	ASP
1	B	322	LEU
1	B	324	TRP

Some 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	B	307	ASN

### 5.3.3 RNA ⓘ

There are no RNA chains 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 ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry

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	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z  > 2$	Counts	RMSZ	# $ Z  > 2$
4	MLI	A	341	-	2,6,6	4.32	1 (50%)	0,7,7	0.00	-
2	PLP	A	350	1	14,15,16	1.33	1 (7%)	20,22,23	1.39	3 (15%)
4	MLI	B	341	-	2,6,6	4.38	2 (100%)	0,7,7	0.00	-
2	PLP	B	350	1	14,15,16	1.09	1 (7%)	20,22,23	1.57	4 (20%)

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

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	341	MLI	C1-C3	5.93	1.53	1.49
4	B	341	MLI	C1-C2	4.57	1.52	1.49
4	B	341	MLI	C1-C3	4.18	1.52	1.49
2	A	350	PLP	C3-C2	-3.54	1.38	1.40
2	B	350	PLP	C3-C2	-3.35	1.38	1.40

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	350	PLP	O2P-P-O1P	3.22	120.98	110.44
2	B	350	PLP	O4P-C5A-C5	3.11	115.58	109.26

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	350	PLP	O4P-C5A-C5	2.97	115.29	109.26
2	A	350	PLP	C2A-C2-N1	2.66	123.77	117.93
2	B	350	PLP	O4P-P-O1P	-2.46	99.49	106.71
2	A	350	PLP	C2A-C2-C3	-2.42	118.07	121.02
2	B	350	PLP	O2P-P-O4P	-2.02	101.06	106.65

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

## 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 ⓘ

### 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	322/339 (94%)	-0.19	1 (0%) 91 93	20, 33, 50, 64	0
1	B	312/339 (92%)	0.57	32 (10%) 7 6	31, 62, 82, 85	0
All	All	634/678 (93%)	0.18	33 (5%) 26 26	20, 43, 80, 85	0

All (33) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	322	LEU	5.3
1	B	138	VAL	4.2
1	B	19	ILE	4.2
1	B	171	LEU	3.8
1	B	7	ILE	3.3
1	B	101	PRO	3.1
1	B	66	GLY	3.0
1	B	61	LEU	3.0
1	B	103	TYR	2.9
1	B	104	ILE	2.9
1	B	296	GLN	2.8
1	B	175	PRO	2.8
1	B	60	ALA	2.8
1	B	63	ALA	2.7
1	B	23	VAL	2.6
1	B	127	TYR	2.5
1	B	116	LEU	2.5
1	B	64	ILE	2.5
1	B	4	GLN	2.4
1	B	168	LEU	2.4
1	B	65	ARG	2.4
1	B	142	ILE	2.4
1	B	306	LYS	2.3
1	B	139	ALA	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	77	LYS	2.3
1	B	204	VAL	2.2
1	B	202	PRO	2.2
1	B	10	ALA	2.2
1	B	195	ILE	2.2
1	B	108	GLN	2.2
1	B	300	THR	2.1
1	B	255	LEU	2.1
1	B	79	VAL	2.1

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

There are no non-standard protein/DNA/RNA residues in this entry.

## 6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 6.4 Ligands ⓘ

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, 95<sup>th</sup> 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	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
4	MLI	A	341	7/7	0.12	0.54	25,30,35,36	0
2	PLP	A	350	15/16	0.12	0.40	20,27,30,30	0
4	MLI	B	341	7/7	0.12	-0.65	44,47,50,51	0
3	MN	B	340	1/1	0.09	-0.74	60,60,60,60	1
2	PLP	B	350	15/16	0.11	-0.91	38,40,42,44	0
3	MN	A	340	1/1	0.07	-1.57	45,45,45,45	1

## 6.5 Other polymers ⓘ

There are no such residues in this entry.