



# Full wwPDB X-ray Structure Validation Report ⓘ

Aug 30, 2023 – 01:35 AM EDT

PDB ID : 3LSQ  
Title : Trypanosoma brucei seryl-tRNA synthetase  
Authors : Larson, E.T.; Merritt, E.A.; Medical Structural Genomics of Pathogenic Protozoa (MSGPP)  
Deposited on : 2010-02-12  
Resolution : 2.10 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](mailto: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 ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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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 : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.35  
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

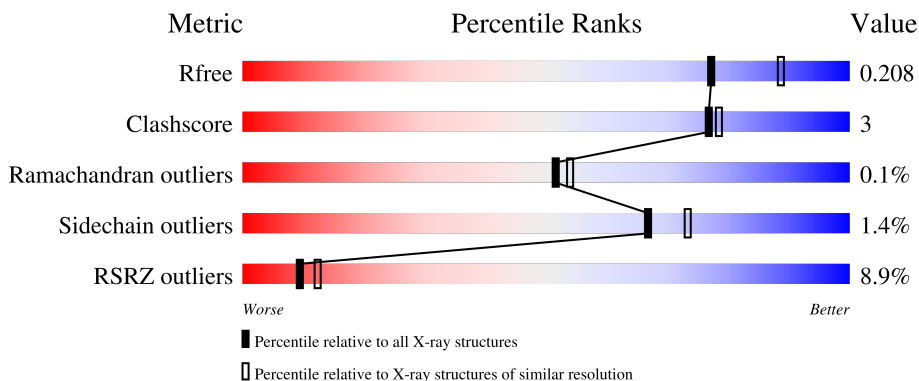
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.10 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))
$R_{free}$	130704	5197 (2.10-2.10)
Clashscore	141614	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (2.10-2.10)
RSRZ outliers	127900	5083 (2.10-2.10)

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  $\geq 3$ , 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  $\leq 5\%$ . 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	A	484	
1	B	484	

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
2	MLI	A	602	-	-	X	-

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7387 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 Seryl-tRNA synthetase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	431	Total	C	N	O	S	0	5	0
			3447	2170	600	648	29			
1	B	458	Total	C	N	O	S	0	6	0
			3649	2293	629	698	29			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-4	GLY	-	expression tag	UNP Q384V4
A	-3	PRO	-	expression tag	UNP Q384V4
A	-2	GLY	-	expression tag	UNP Q384V4
A	-1	SER	-	expression tag	UNP Q384V4
A	0	MET	-	expression tag	UNP Q384V4
B	-4	GLY	-	expression tag	UNP Q384V4
B	-3	PRO	-	expression tag	UNP Q384V4
B	-2	GLY	-	expression tag	UNP Q384V4
B	-1	SER	-	expression tag	UNP Q384V4
B	0	MET	-	expression tag	UNP Q384V4

- Molecule 2 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
2	A	1	Total C O 7 3 4	0	0

- Molecule 3 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total Na 1 1	0	0

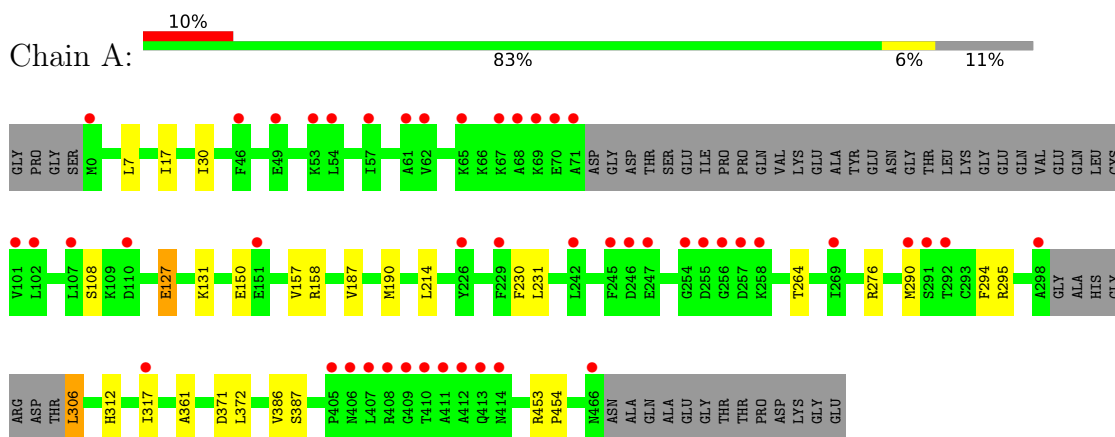
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	142	Total O 142 142	0	0
4	B	141	Total O 141 141	0	0

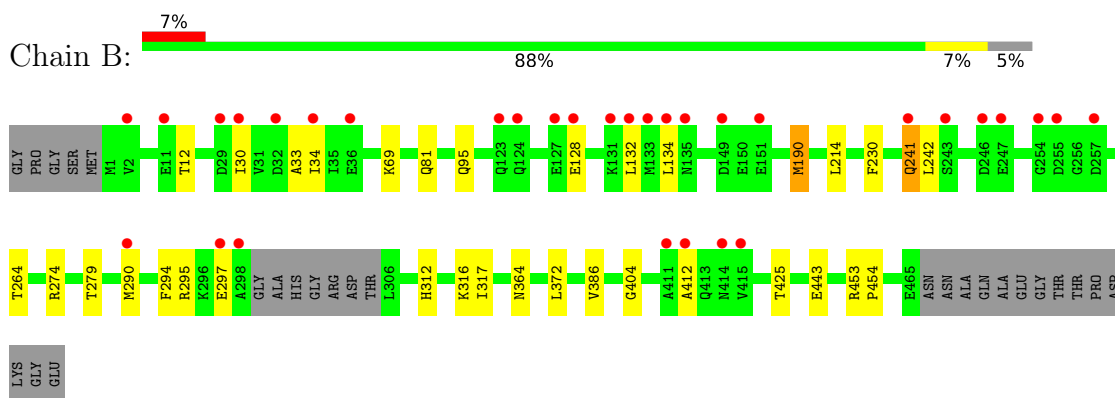
### 3 Residue-property plots [i](#)

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: Seryl-tRNA synthetase



- Molecule 1: Seryl-tRNA synthetase



## 4 Data and refinement statistics

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	176.02Å 176.02Å 248.26Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	37.45 – 2.10 36.43 – 2.10	Depositor EDS
% Data completeness (in resolution range)	99.5 (37.45-2.10) 99.5 (36.43-2.10)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.98 (at 2.10Å)	Xtrriage
Refinement program	REFMAC refmac_5.5.0106	Depositor
R, $R_{free}$	0.171 , 0.203 0.179 , 0.208	Depositor DCC
$R_{free}$ test set	4328 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	31.5	Xtrriage
Anisotropy	0.286	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.38 , 48.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	7387	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	43.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.29% of the height of the origin peak. No significant pseudotranslation is detected.*

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

<sup>2</sup>Theoretical 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: NA, 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	Bond lengths		Bond angles	
		RMSZ	# $ Z  > 5$	RMSZ	# $ Z  > 5$
1	A	0.69	0/3521	0.71	2/4737 (0.0%)
1	B	0.68	1/3729 (0.0%)	0.70	2/5024 (0.0%)
All	All	0.69	1/7250 (0.0%)	0.71	4/9761 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	443	GLU	CB-CG	5.04	1.61	1.52

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	230	PHE	CB-CA-C	-6.68	97.04	110.40
1	A	230	PHE	CB-CA-C	-6.17	98.06	110.40
1	B	190	MET	CG-SD-CE	5.45	108.92	100.20
1	A	295	ARG	NE-CZ-NH2	-5.01	117.79	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	3447	0	3454	18	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	3649	0	3638	26	0
2	A	7	0	2	2	0
3	A	1	0	0	0	0
4	A	142	0	0	1	0
4	B	141	0	0	0	0
All	All	7387	0	7094	42	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 (42) 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:B:30:ILE:HG22	1:B:34:ILE:CD1	1.97	0.95
1:B:30:ILE:HG22	1:B:34:ILE:HD11	1.50	0.91
1:B:214:LEU:HD13	1:B:290[B]:MET:SD	2.14	0.87
1:B:134:LEU:HD11	1:B:242:LEU:HD12	1.60	0.82
1:B:33:ALA:HB1	1:B:132:LEU:HD23	1.62	0.81
1:B:30:ILE:O	1:B:34:ILE:HD12	1.83	0.79
1:B:290[B]:MET:SD	1:B:317:ILE:CD1	2.74	0.76
1:B:30:ILE:CG2	1:B:34:ILE:HD11	2.14	0.75
1:A:214:LEU:HD13	1:A:290[B]:MET:SD	2.30	0.72
1:B:290[B]:MET:SD	1:B:317:ILE:HD12	2.39	0.61
1:A:127[A]:GLU:HA	1:A:127[A]:GLU:OE1	2.01	0.60
1:A:276:ARG:HD3	2:A:602:MLI:H11	1.85	0.58
1:B:290[B]:MET:SD	1:B:317:ILE:HD13	2.44	0.58
1:B:33:ALA:HB1	1:B:132:LEU:CD2	2.34	0.57
1:A:317:ILE:HD12	1:A:317:ILE:N	2.20	0.56
1:B:128:GLU:O	1:B:132:LEU:HD13	2.04	0.56
1:A:231:LEU:HD22	1:B:190:MET:HG2	1.89	0.53
1:B:241:GLN:HE22	1:B:364:ASN:ND2	2.07	0.52
1:A:306:LEU:HD23	4:A:942:HOH:O	2.07	0.52
1:B:372:LEU:HB3	1:B:386:VAL:HB	1.93	0.51
1:B:81:GLN:OE1	1:B:95:GLN:NE2	2.42	0.51
1:B:30:ILE:CG2	1:B:34:ILE:CD1	2.78	0.51
1:B:33:ALA:CB	1:B:132:LEU:HD23	2.35	0.51
1:A:372:LEU:HB3	1:A:386:VAL:HB	1.94	0.50
1:A:187:VAL:HA	1:A:190:MET:HE2	1.95	0.48
1:A:371:ASP:OD1	1:A:387:SER:HB3	2.14	0.48
1:A:157:VAL:HG12	1:A:158:ARG:HG3	1.96	0.47
1:B:295:ARG:NH1	1:B:297:GLU:OE2	2.45	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:7:LEU:HD23	1:A:17:ILE:CG1	2.47	0.45
1:A:150:GLU:HG3	1:A:361:ALA:CB	2.48	0.44
1:A:231:LEU:CD2	1:B:190:MET:HG2	2.48	0.43
1:B:294:PHE:HA	1:B:312:HIS:O	2.18	0.43
1:B:316:LYS:HD2	1:B:425:THR:HG23	2.00	0.43
1:B:30:ILE:HG22	1:B:34:ILE:HD12	1.91	0.43
1:A:187:VAL:HA	1:A:190:MET:CE	2.49	0.43
1:A:294:PHE:HA	1:A:312:HIS:O	2.19	0.42
1:B:134:LEU:CD1	1:B:242:LEU:HD12	2.41	0.42
1:A:187:VAL:O	1:A:190:MET:HB2	2.20	0.41
1:A:276:ARG:HH11	2:A:602:MLI:H11	1.85	0.41
1:B:453:ARG:HB2	1:B:454:PRO:HD3	2.02	0.41
1:A:453:ARG:N	1:A:454:PRO:CD	2.84	0.40
1:B:279:THR:HA	1:B:404:GLY:O	2.21	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	430/484 (89%)	425 (99%)	5 (1%)	0	100	100
1	B	460/484 (95%)	451 (98%)	8 (2%)	1 (0%)	47	49
All	All	890/968 (92%)	876 (98%)	13 (2%)	1 (0%)	51	54

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	412	ALA

### 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	377/412 (92%)	370 (98%)	7 (2%)	57	63
1	B	400/412 (97%)	395 (99%)	5 (1%)	69	75
All	All	777/824 (94%)	765 (98%)	12 (2%)	67	71

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	30	ILE
1	A	108	SER
1	A	127[A]	GLU
1	A	127[B]	GLU
1	A	131	LYS
1	A	264	THR
1	A	306	LEU
1	B	12	THR
1	B	69	LYS
1	B	241	GLN
1	B	264	THR
1	B	274	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	241	GLN

### 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 2 ligands modelled in this entry, 1 is monoatomic - leaving 1 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
2	MLI	A	602	-	6,6,6	1.88	2 (33%)	7,7,7	1.21	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
2	MLI	A	602	-	-	2/4/4/4	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	602	MLI	C1-C2	2.86	1.55	1.51
2	A	602	MLI	C1-C3	2.54	1.55	1.51

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	602	MLI	C2-C1-C3-O8
2	A	602	MLI	C2-C1-C3-O9

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	602	MLI	2	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<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	431/484 (89%)	0.39	47 (10%) <b>5</b> <b>7</b>	20, 33, 93, 129	0
1	B	458/484 (94%)	0.13	32 (6%) <b>16</b> <b>20</b>	21, 36, 83, 111	0
All	All	889/968 (91%)	0.26	79 (8%) <b>9</b> <b>12</b>	20, 35, 86, 129	0

All (79) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	132	LEU	10.9
1	A	298	ALA	8.7
1	A	255	ASP	8.3
1	B	298	ALA	8.2
1	B	414	ASN	7.1
1	A	408	ARG	6.8
1	A	410	THR	6.8
1	A	54	LEU	6.3
1	A	69	LYS	5.7
1	A	254	GLY	5.6
1	A	68	ALA	5.6
1	B	255	ASP	5.5
1	A	57	ILE	5.2
1	B	30	ILE	4.7
1	A	246	ASP	4.5
1	A	407	LEU	4.5
1	A	406	ASN	4.4
1	A	46	PHE	4.4
1	B	246	ASP	4.3
1	B	131	LYS	4.3
1	B	415	VAL	4.3
1	A	409	GLY	4.1
1	A	414	ASN	4.0
1	A	67	LYS	3.9

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
1	A	102	LEU	3.9
1	B	151	GLU	3.7
1	A	405	PRO	3.7
1	B	123	GLN	3.7
1	A	71	ALA	3.7
1	B	134	LEU	3.6
1	A	290[A]	MET	3.6
1	B	257	ASP	3.5
1	A	413	GLN	3.4
1	A	247	GLU	3.4
1	A	257	ASP	3.3
1	A	70	GLU	3.3
1	A	101	VAL	3.3
1	A	242	LEU	3.3
1	B	254	GLY	3.2
1	A	61	ALA	3.2
1	B	133[A]	MET	3.2
1	B	412	ALA	3.2
1	A	0	MET	3.1
1	B	243	SER	3.1
1	B	247	GLU	3.1
1	B	127	GLU	3.0
1	A	107	LEU	3.0
1	B	34	ILE	3.0
1	B	32	ASP	2.9
1	A	245	PHE	2.9
1	A	411	ALA	2.8
1	B	29	ASP	2.8
1	A	151	GLU	2.8
1	B	11	GLU	2.7
1	B	241	GLN	2.7
1	A	53	LYS	2.7
1	B	36	GLU	2.6
1	B	124	GLN	2.6
1	B	135	ASN	2.6
1	A	412	ALA	2.5
1	B	290[A]	MET	2.5
1	B	128	GLU	2.5
1	A	65	LYS	2.5
1	A	49	GLU	2.5
1	A	466	ASN	2.5
1	B	411	ALA	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	62	VAL	2.4
1	B	149	ASP	2.4
1	A	258	LYS	2.3
1	A	226	TYR	2.3
1	A	269	ILE	2.3
1	A	110	ASP	2.2
1	A	256	GLY	2.1
1	B	297	GLU	2.1
1	A	317	ILE	2.1
1	A	292	THR	2.1
1	A	229	PHE	2.1
1	A	291	SER	2.0
1	B	2	VAL	2.0

## 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 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<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	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
2	MLI	A	602	7/7	0.87	0.19	40,51,62,63	0
3	NA	A	600	1/1	0.96	0.10	25,25,25,25	1

## 6.5 Other polymers [i](#)

There are no such residues in this entry.