



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 02:29 AM UTC

PDB ID : 6OR9 / pdb_00006or9
Title : Structure of L-lactate dehydrogenase from Trichoplusia ni
Authors : Seattle Structural Genomics Center for Infectious Disease (SSGCID)
Deposited on : 2019-04-29
Resolution : 1.80 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 ⓘ 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 ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

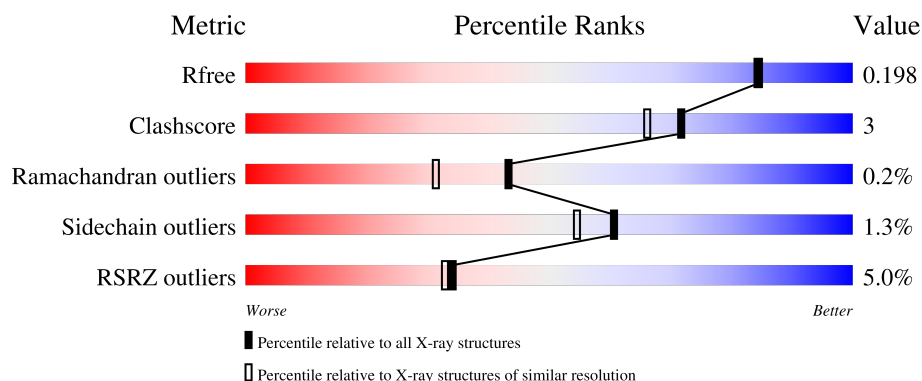
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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}	180053	7662 (1.80-1.80)
Clashscore	190562	8479 (1.80-1.80)
Ramachandran outliers	187476	8391 (1.80-1.80)
Sidechain outliers	187428	8390 (1.80-1.80)
RSRZ outliers	180081	7663 (1.80-1.80)

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 $\leq 5\%$. 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	364	4% 79% 16%
1	B	364	2% 79% 15%
1	C	364	5% 78% 16%
1	D	364	6% 73% 18%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 10401 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 L-lactate dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	306	Total	C	N	O	S	0	12	0
			2330	1471	408	438	13			
1	B	308	Total	C	N	O	S	0	19	0
			2398	1519	414	451	14			
1	C	305	Total	C	N	O	S	0	9	0
			2306	1461	397	435	13			
1	D	299	Total	C	N	O	S	0	7	0
			2241	1420	380	428	13			

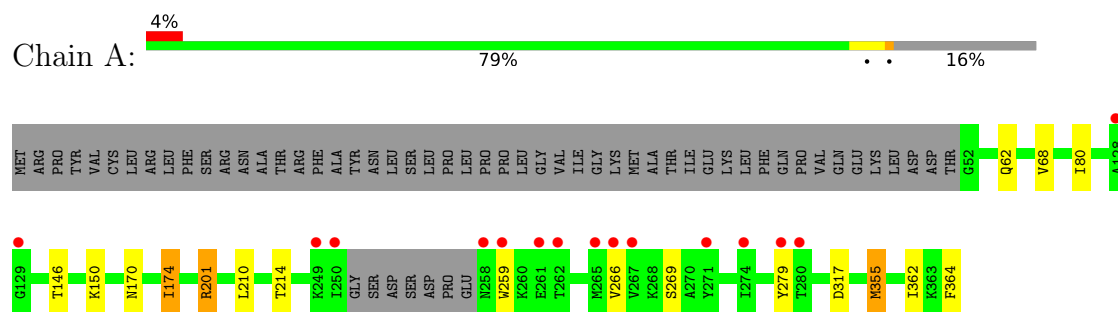
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	287	Total	O	0	6
			293	293		
2	B	322	Total	O	0	13
			334	334		
2	C	276	Total	O	0	8
			284	284		
2	D	212	Total	O	0	4
			215	215		

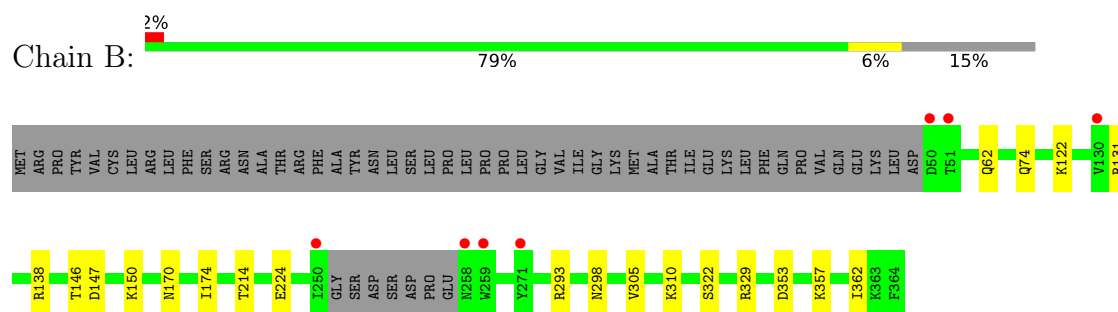
3 Residue-property plots

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.

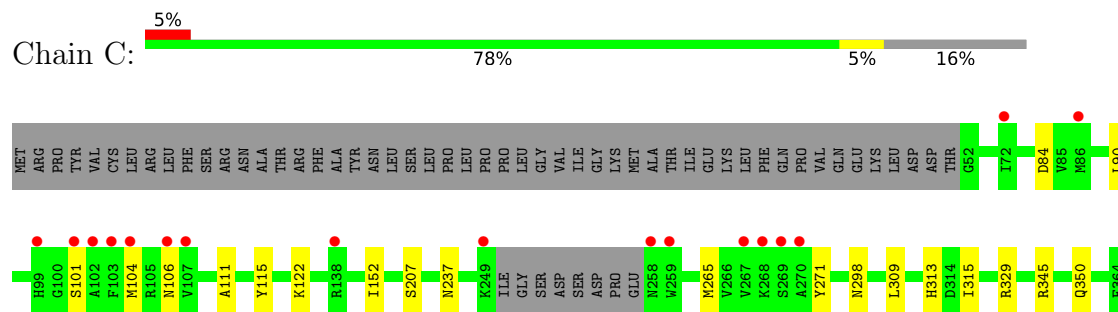
- Molecule 1: L-lactate dehydrogenase



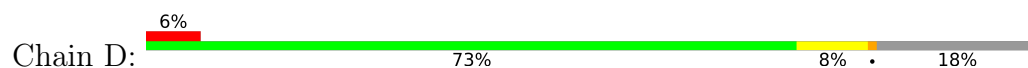
- Molecule 1: L-lactate dehydrogenase

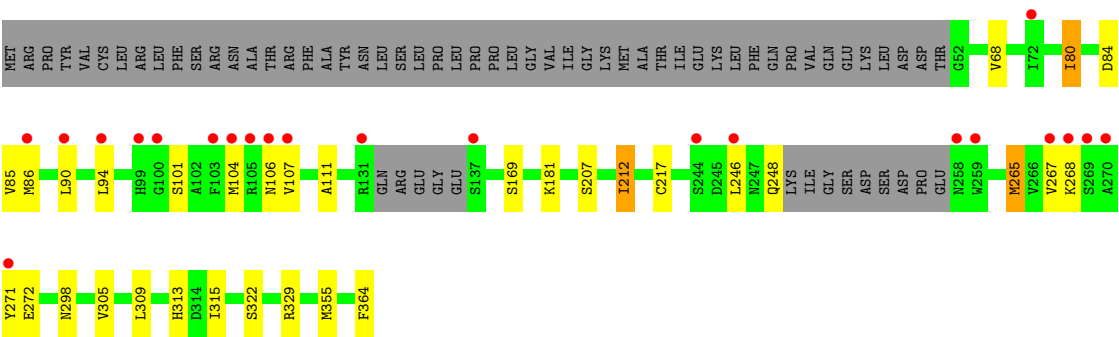


- Molecule 1: L-lactate dehydrogenase



- Molecule 1: L-lactate dehydrogenase





4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	63.39Å 76.48Å 79.52Å 113.06° 109.84° 88.70°	Depositor
Resolution (Å)	48.91 – 1.80 48.91 – 1.80	Depositor EDS
% Data completeness (in resolution range)	97.0 (48.91-1.80) 97.0 (48.91-1.80)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.73 (at 1.79Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.158 , 0.195 0.160 , 0.198	Depositor DCC
R_{free} test set	1971 reflections (1.66%)	wwPDB-VP
Wilson B-factor (Å ²)	22.1	Xtriage
Anisotropy	0.364	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 53.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	10401	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.78% 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](#)

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	0.34	0/2399	0.51	0/3254
1	B	0.36	0/2487	0.54	0/3373
1	C	0.34	0/2367	0.53	0/3215
1	D	0.29	0/2295	0.49	0/3121
All	All	0.33	0/9548	0.52	0/12963

There are no bond length outliers.

There are no bond angle outliers.

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	2330	0	2385	15	0
1	B	2398	0	2465	16	0
1	C	2306	0	2348	14	0
1	D	2241	0	2252	17	0
2	A	293	0	0	4	0
2	B	334	0	0	7	0
2	C	284	0	0	5	0
2	D	215	0	0	1	0
All	All	10401	0	9450	62	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 (62) 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:224:GLU:OE1	2:B:401:HOH:O	2.02	0.77
1:B:122[A]:LYS:NZ	2:B:403[A]:HOH:O	2.22	0.72
1:C:350:GLN:OE1	2:C:401:HOH:O	2.08	0.70
1:A:201[B]:ARG:NH2	1:A:266:VAL:O	2.25	0.70
1:C:101:SER:HA	1:C:104:MET:HE2	1.74	0.69
1:A:317:ASP:OD1	2:A:403:HOH:O	2.12	0.66
1:A:364:PHE:OXT	2:A:402:HOH:O	2.12	0.66
1:D:268:LYS:O	1:D:272[A]:GLU:HG2	1.98	0.62
1:B:138:ARG:NH2	2:B:408:HOH:O	2.33	0.61
1:B:353[B]:ASP:OD1	1:B:357:LYS:NZ	2.34	0.61
1:A:68:VAL:HG13	1:A:80[A]:ILE:HD13	1.83	0.60
1:C:309:LEU:HD21	1:C:315[B]:ILE:HD13	1.84	0.59
1:C:298:ASN:HB2	1:C:329:ARG:HG2	1.85	0.58
1:D:101:SER:HA	1:D:104:MET:HE2	1.85	0.57
1:A:170:ASN:ND2	2:A:407:HOH:O	2.36	0.57
1:B:174[A]:ILE:HD11	1:B:362:ILE:HD11	1.89	0.55
1:D:84:ASP:H	1:D:90:LEU:HD11	1.73	0.54
1:C:237:ASN:OD1	2:C:402:HOH:O	2.19	0.53
1:D:246:LEU:HD22	1:D:248:GLN:HG2	1.90	0.52
1:B:122[B]:LYS:HE2	1:B:329:ARG:NE	2.25	0.51
1:B:298:ASN:HB2	1:B:329:ARG:HG2	1.91	0.51
1:C:104:MET:O	2:C:403:HOH:O	2.19	0.50
1:B:62[A]:GLN:NE2	1:B:131:ARG:HH22	2.09	0.50
1:C:90:LEU:HD22	1:C:111:ALA:HB1	1.95	0.49
1:A:201[A]:ARG:HG2	1:A:266:VAL:HG22	1.95	0.48
1:C:345:ARG:HD3	2:C:489:HOH:O	2.13	0.48
1:A:210:LEU:HD21	1:A:259:TRP:HH2	1.78	0.48
1:D:298:ASN:HB2	1:D:329:ARG:HG2	1.95	0.48
1:B:170:ASN:ND2	2:B:402:HOH:O	2.13	0.47
1:C:313:HIS:HB2	1:C:315[B]:ILE:HD12	1.96	0.47
1:B:146:THR:HG22	1:B:150:LYS:HE2	1.97	0.47
1:B:147:ASP:HA	1:B:150:LYS:HE3	1.95	0.47
1:B:214[B]:THR:HG21	2:B:426:HOH:O	2.15	0.47
1:D:68:VAL:HG13	1:D:80:ILE:HD12	1.97	0.46
1:D:305:VAL:O	1:D:322:SER:HA	2.15	0.46
1:C:115:TYR:OH	1:C:152:ILE:HG12	2.15	0.46
1:A:170:ASN:ND2	2:A:401:HOH:O	2.07	0.46
1:A:146:THR:HG22	1:A:150:LYS:HE3	1.98	0.45
1:A:174[A]:ILE:HD11	1:A:362:ILE:HD11	1.99	0.45
1:D:104:MET:HE3	1:D:107:VAL:HG21	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:170:ASN:ND2	2:B:416[A]:HOH:O	2.47	0.44
1:B:74:GLN:O	1:B:293:ARG:NH2	2.50	0.44
1:C:101:SER:HA	1:C:104:MET:CE	2.45	0.44
1:D:265[A]:MET:HE2	1:D:265[A]:MET:HB2	1.88	0.44
1:D:90:LEU:HD22	1:D:111:ALA:HB1	1.99	0.44
1:C:84:ASP:H	1:C:90:LEU:HD11	1.83	0.43
1:D:212[A]:ILE:HD11	1:D:217:CYS:SG	2.58	0.43
1:A:174[B]:ILE:HD11	1:A:355:MET:HG3	2.01	0.43
1:B:310:LYS:NZ	2:B:419:HOH:O	2.50	0.43
1:A:214[B]:THR:HG21	2:D:406:HOH:O	2.18	0.43
1:D:84:ASP:OD2	1:D:85:VAL:N	2.46	0.42
1:D:181:LYS:HD3	1:D:364:PHE:CZ	2.54	0.42
1:D:309:LEU:HD21	1:D:315[B]:ILE:HD13	2.01	0.42
1:C:265:MET:HE2	1:C:265:MET:HB2	1.90	0.42
1:A:201[B]:ARG:HH11	1:A:201[B]:ARG:HG3	1.85	0.42
1:D:313:HIS:HB2	1:D:315[B]:ILE:HD12	2.03	0.41
1:C:122:LYS:HD2	2:C:411:HOH:O	2.20	0.41
1:B:305:VAL:O	1:B:322:SER:HA	2.20	0.41
1:A:62[A]:GLN:HG2	1:A:279:TYR:CE1	2.56	0.41
1:D:94:LEU:HB2	1:D:111:ALA:HB3	2.03	0.41
1:A:201[B]:ARG:NH2	1:A:269:SER:OG	2.54	0.40
1:D:85:VAL:HG23	1:D:86:MET:HG3	2.04	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	314/364 (86%)	309 (98%)	5 (2%)	0	100	100
1	B	323/364 (89%)	318 (98%)	5 (2%)	0	100	100
1	C	310/364 (85%)	302 (97%)	7 (2%)	1 (0%)	36	25

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	D	300/364 (82%)	294 (98%)	5 (2%)	1 (0%)	36	25
All	All	1247/1456 (86%)	1223 (98%)	22 (2%)	2 (0%)	43	31

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	106	ASN
1	C	106	ASN

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	260/312 (83%)	255 (98%)	5 (2%)	50	41
1	B	271/312 (87%)	271 (100%)	0	100	100
1	C	256/312 (82%)	254 (99%)	2 (1%)	73	70
1	D	247/312 (79%)	237 (96%)	10 (4%)	28	15
All	All	1034/1248 (83%)	1017 (98%)	17 (2%)	61	47

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	174[A]	ILE
1	A	174[B]	ILE
1	A	201[A]	ARG
1	A	201[B]	ARG
1	A	355	MET
1	C	207	SER
1	C	271	TYR
1	D	80	ILE
1	D	169	SER
1	D	207	SER
1	D	212[A]	ILE
1	D	212[B]	ILE

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Mol	Chain	Res	Type
1	D	265[A]	MET
1	D	265[B]	MET
1	D	267	VAL
1	D	271	TYR
1	D	355	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	196	ASN
1	A	218	HIS
1	A	290	GLN
1	B	196	ASN
1	B	338	GLN
1	B	347	GLN
1	C	99	HIS
1	C	188	HIS
1	C	196	ASN
1	C	218	HIS
1	C	290	GLN
1	D	196	ASN
1	D	218	HIS
1	D	290	GLN
1	D	350	GLN

5.3.3 RNA ⓘ

There are no RNA molecules 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 oligosaccharides in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

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	A	306/364 (84%)	-0.24	15 (4%) 35 33	10, 23, 51, 74	12 (3%)
1	B	308/364 (84%)	-0.43	7 (2%) 61 61	9, 21, 41, 80	19 (6%)
1	C	305/364 (83%)	-0.16	17 (5%) 30 29	14, 25, 49, 79	9 (2%)
1	D	299/364 (82%)	0.11	22 (7%) 20 19	13, 29, 57, 82	7 (2%)
All	All	1218/1456 (83%)	-0.18	61 (5%) 34 33	9, 24, 52, 82	47 (3%)

All (61) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	259	TRP	5.3
1	B	258	ASN	4.7
1	A	250	ILE	4.2
1	D	270	ALA	4.1
1	D	99	HIS	4.1
1	C	270	ALA	3.9
1	D	72	ILE	3.8
1	D	269	SER	3.8
1	C	259	TRP	3.8
1	D	90	LEU	3.6
1	A	258	ASN	3.5
1	D	106	ASN	3.4
1	B	250	ILE	3.4
1	D	131	ARG	3.3
1	D	103	PHE	3.2
1	D	267	VAL	3.2
1	D	100	GLY	3.1
1	D	105	ARG	3.1
1	C	269	SER	3.1
1	D	137	SER	3.0
1	A	280	THR	3.0

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Mol	Chain	Res	Type	RSRZ
1	C	249	LYS	2.9
1	D	86	MET	2.9
1	C	267	VAL	2.8
1	A	128	ALA	2.8
1	C	107	VAL	2.8
1	C	104	MET	2.8
1	A	271	TYR	2.8
1	D	271	TYR	2.7
1	D	258	ASN	2.7
1	C	103	PHE	2.7
1	A	129	GLY	2.7
1	A	262	THR	2.7
1	B	259	TRP	2.6
1	D	246	LEU	2.6
1	A	266	VAL	2.6
1	C	268	LYS	2.6
1	A	274	ILE	2.5
1	A	279	TYR	2.5
1	C	106	ASN	2.5
1	C	101	SER	2.5
1	D	104	MET	2.5
1	D	107	VAL	2.5
1	D	268	LYS	2.4
1	C	258	ASN	2.4
1	B	130	VAL	2.4
1	D	244	SER	2.3
1	B	51	THR	2.2
1	D	259	TRP	2.2
1	A	249	LYS	2.2
1	A	261	GLU	2.2
1	C	102	ALA	2.2
1	A	265	MET	2.1
1	C	86	MET	2.1
1	D	94	LEU	2.1
1	C	99	HIS	2.1
1	B	50	ASP	2.1
1	C	138	ARG	2.1
1	A	267	VAL	2.1
1	C	72	ILE	2.0
1	B	271	TYR	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.