



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 06:31 AM UTC

PDB ID : 3NVT / pdb_00003nvt
Title : 1.95 Angstrom crystal structure of a bifunctional 3-deoxy-7-phosphoheptulon
ate synthase/chorismate mutase (aroA) from *Listeria monocytogenes* EGD-e
Authors : Halavaty, A.S.; Light, S.H.; Minasov, G.; Shuvalova, L.; Kwon, K.; Anderson,
W.F.; Center for Structural Genomics of Infectious Diseases (CSGID)
Deposited on : 2010-07-08
Resolution : 1.95 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
Mogul	:	2022.3.0, CSD as543be (2022)
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

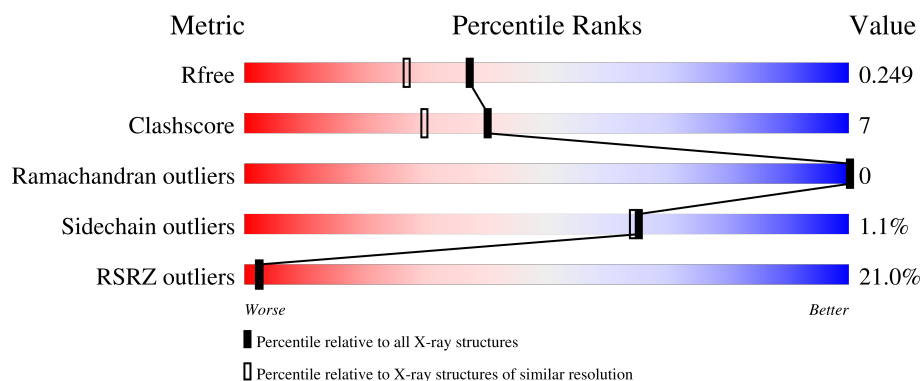
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.95 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	3494 (1.96-1.96)
Clashscore	190562	3612 (1.96-1.96)
Ramachandran outliers	187476	3587 (1.96-1.96)
Sidechain outliers	187428	3587 (1.96-1.96)
RSRZ outliers	180081	3495 (1.96-1.96)

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	385	19% 78% 11% 10%
1	B	385	17% 77% 8% 15%

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
3	ACT	A	363	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6105 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 3-deoxy-D-arabino-heptulosonate 7-phosphate synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	345	Total	C	N	O	S	0	13	0
			2768	1751	471	536	10			
1	B	327	Total	C	N	O	S	0	7	0
			2562	1628	431	493	10			

There are 48 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-23	MET	-	expression tag	UNP Q8Y6T2
A	-22	HIS	-	expression tag	UNP Q8Y6T2
A	-21	HIS	-	expression tag	UNP Q8Y6T2
A	-20	HIS	-	expression tag	UNP Q8Y6T2
A	-19	HIS	-	expression tag	UNP Q8Y6T2
A	-18	HIS	-	expression tag	UNP Q8Y6T2
A	-17	HIS	-	expression tag	UNP Q8Y6T2
A	-16	SER	-	expression tag	UNP Q8Y6T2
A	-15	SER	-	expression tag	UNP Q8Y6T2
A	-14	GLY	-	expression tag	UNP Q8Y6T2
A	-13	VAL	-	expression tag	UNP Q8Y6T2
A	-12	ASP	-	expression tag	UNP Q8Y6T2
A	-11	LEU	-	expression tag	UNP Q8Y6T2
A	-10	GLY	-	expression tag	UNP Q8Y6T2
A	-9	THR	-	expression tag	UNP Q8Y6T2
A	-8	GLU	-	expression tag	UNP Q8Y6T2
A	-7	ASN	-	expression tag	UNP Q8Y6T2
A	-6	LEU	-	expression tag	UNP Q8Y6T2
A	-5	TYR	-	expression tag	UNP Q8Y6T2
A	-4	PHE	-	expression tag	UNP Q8Y6T2
A	-3	GLN	-	expression tag	UNP Q8Y6T2
A	-2	SER	-	expression tag	UNP Q8Y6T2
A	-1	ASN	-	expression tag	UNP Q8Y6T2
A	0	ALA	-	expression tag	UNP Q8Y6T2
B	-23	MET	-	expression tag	UNP Q8Y6T2

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-22	HIS	-	expression tag	UNP Q8Y6T2
B	-21	HIS	-	expression tag	UNP Q8Y6T2
B	-20	HIS	-	expression tag	UNP Q8Y6T2
B	-19	HIS	-	expression tag	UNP Q8Y6T2
B	-18	HIS	-	expression tag	UNP Q8Y6T2
B	-17	HIS	-	expression tag	UNP Q8Y6T2
B	-16	SER	-	expression tag	UNP Q8Y6T2
B	-15	SER	-	expression tag	UNP Q8Y6T2
B	-14	GLY	-	expression tag	UNP Q8Y6T2
B	-13	VAL	-	expression tag	UNP Q8Y6T2
B	-12	ASP	-	expression tag	UNP Q8Y6T2
B	-11	LEU	-	expression tag	UNP Q8Y6T2
B	-10	GLY	-	expression tag	UNP Q8Y6T2
B	-9	THR	-	expression tag	UNP Q8Y6T2
B	-8	GLU	-	expression tag	UNP Q8Y6T2
B	-7	ASN	-	expression tag	UNP Q8Y6T2
B	-6	LEU	-	expression tag	UNP Q8Y6T2
B	-5	TYR	-	expression tag	UNP Q8Y6T2
B	-4	PHE	-	expression tag	UNP Q8Y6T2
B	-3	GLN	-	expression tag	UNP Q8Y6T2
B	-2	SER	-	expression tag	UNP Q8Y6T2
B	-1	ASN	-	expression tag	UNP Q8Y6T2
B	0	ALA	-	expression tag	UNP Q8Y6T2

- Molecule 2 is MANGANESE (II) ION (CCD ID: MN) (formula: Mn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Mn 2 2	0	1
2	B	1	Total Mn 2 2	0	1

- Molecule 3 is ACETATE ION (CCD ID: ACT) (formula: C₂H₃O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			4	2	2		
3	B	1	Total	C	O	0	0
			4	2	2		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	394	Total	O	0	14
			399	399		
4	B	362	Total	O	0	5
			364	364		

- Molecule 1: 3-deoxy-D-arabino-heptulosonate 7-phosphate synthase



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	111.57Å 111.79Å 81.05Å 90.00° 127.63° 90.00°	Depositor
Resolution (Å)	29.04 – 1.95 29.04 – 1.95	Depositor EDS
% Data completeness (in resolution range)	99.5 (29.04-1.95) 99.7 (29.04-1.95)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.58 (at 1.95Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.154 , 0.198 0.214 , 0.249	Depositor DCC
R_{free} test set	2907 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	19.5	Xtriage
Anisotropy	0.270	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 40.4	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	0.94	EDS
Total number of atoms	6105	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.89% 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](#)

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

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	0.58	0/2806	0.82	2/3786 (0.1%)
1	B	0.59	0/2595	0.81	1/3504 (0.0%)
All	All	0.59	0/5401	0.82	3/7290 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	190	VAL	N-CA-C	5.79	118.45	112.90
1	B	190	VAL	N-CA-C	5.49	118.17	112.90
1	A	11	THR	N-CA-C	-5.05	106.28	112.90

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	2768	0	2822	47	0
1	B	2562	0	2636	31	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
3	A	4	0	3	4	0
3	B	4	0	3	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	399	0	0	18	0
4	B	364	0	0	9	0
All	All	6105	0	5464	71	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 7.

All (71) 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:A:320[B]:MET:SD	4:A:683:HOH:O	2.23	0.97
1:B:320[B]:MET:SD	4:B:499:HOH:O	2.38	0.81
1:A:285[B]:GLU:HG2	4:A:528:HOH:O	1.88	0.72
1:B:150[A]:ARG:NH1	4:B:687[A]:HOH:O	2.21	0.72
1:B:54:MET:HE3	1:B:79:PHE:CZ	2.24	0.72
1:B:150[A]:ARG:HH11	1:B:150[A]:ARG:HB3	1.58	0.69
1:A:150[B]:ARG:CD	4:A:649:HOH:O	2.41	0.67
1:B:356:VAL:HG22	4:B:551:HOH:O	1.95	0.66
1:A:99:LYS:NZ	4:A:413:HOH:O	2.30	0.64
1:B:150[A]:ARG:NH2	1:B:322:GLU:OE2	2.30	0.64
1:A:150[B]:ARG:HD2	4:A:649:HOH:O	1.97	0.64
1:A:291:MET:HE3	1:A:318:GLY:HA3	1.82	0.61
1:A:23:LEU:HD13	1:B:20:LEU:HD23	1.83	0.60
1:A:298:THR:HG21	4:A:707:HOH:O	2.00	0.60
1:A:155:LYS:NZ	3:A:363:ACT:H1	2.17	0.59
1:A:267:GLU:HG2	4:A:621:HOH:O	2.02	0.59
1:A:81:ALA:O	1:A:84:GLU:HG2	2.02	0.58
1:A:155:LYS:HZ2	3:A:363:ACT:H1	1.67	0.58
1:B:7:GLU:O	1:B:11:THR:HG23	2.04	0.58
1:B:126:CYS:HA	1:B:150[A]:ARG:HH22	1.70	0.57
1:A:98:ARG:O	1:A:102:LYS:HD3	2.06	0.56
1:A:108:THR:HG22	4:A:635:HOH:O	2.07	0.55
1:B:285:GLU:HG2	4:B:564:HOH:O	2.06	0.55
1:A:92:LYS:NZ	1:B:69:SER:OG	2.39	0.54
1:A:320[B]:MET:HG3	4:A:649:HOH:O	2.07	0.54
1:A:150[A]:ARG:NH2	1:A:322:GLU:OE2	2.41	0.53
1:A:110:LYS:HG3	1:A:184:GLY:HA3	1.91	0.53
1:A:110:LYS:HG3	1:A:184:GLY:CA	2.38	0.53
1:A:23:LEU:HD22	1:B:20:LEU:HG	1.90	0.53
1:A:150[A]:ARG:HD2	4:A:619[A]:HOH:O	2.10	0.52
1:A:298:THR:CG2	4:A:707:HOH:O	2.57	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:231:LYS:NZ	3:A:363:ACT:O	2.32	0.51
1:B:150[A]:ARG:NH1	1:B:150[A]:ARG:HB3	2.24	0.51
1:A:62:ASN:HA	1:B:21:LEU:HD21	1.93	0.50
1:A:20:LEU:HD23	1:B:23:LEU:HD23	1.94	0.50
1:A:150[B]:ARG:O	1:A:150[B]:ARG:HG3	2.13	0.48
1:B:150[A]:ARG:HD2	4:B:683[A]:HOH:O	2.12	0.48
1:B:255:LYS:NZ	4:B:636:HOH:O	2.45	0.48
3:A:363:ACT:H3	4:A:697:HOH:O	2.14	0.48
1:A:150[B]:ARG:NH1	4:A:683:HOH:O	2.46	0.48
1:B:186:ILE:HG12	1:B:204:VAL:HB	1.96	0.48
1:A:178[B]:SER:HB2	4:A:746[B]:HOH:O	2.14	0.47
1:A:150[A]:ARG:CD	4:A:619[A]:HOH:O	2.63	0.46
1:A:45:ARG:HD3	1:A:46:PHE:H	1.80	0.45
1:A:38:LYS:HG2	1:B:6:LEU:HD11	1.98	0.45
1:A:124:GLY:HA3	1:A:150[B]:ARG:HG3	1.96	0.45
1:A:150[A]:ARG:HH11	1:A:150[A]:ARG:HB3	1.81	0.45
1:B:126:CYS:HA	1:B:150[A]:ARG:NH2	2.31	0.45
1:B:255:LYS:NZ	4:B:658:HOH:O	2.50	0.45
1:A:261:ARG:HD3	1:A:261:ARG:C	2.42	0.45
1:A:251:GLN:HG2	4:A:565:HOH:O	2.16	0.44
1:A:69:SER:O	1:A:73:LYS:HG3	2.18	0.44
1:B:14:ASP:OD2	4:B:720:HOH:O	2.21	0.44
1:B:324:HIS:O	1:B:339:ASP:HA	2.17	0.44
1:A:67[A]:GLU:HG2	4:A:727:HOH:O	2.18	0.44
1:B:54:MET:HE3	1:B:79:PHE:CE2	2.52	0.43
1:A:21:LEU:HD21	1:B:62:ASN:HA	1.99	0.43
1:A:186:ILE:HG12	1:A:204:VAL:HB	2.01	0.43
1:A:268:LYS:HG2	4:A:609:HOH:O	2.18	0.43
1:A:305:LEU:HB3	1:A:306:PRO:HD3	2.01	0.43
1:B:305:LEU:HB3	1:B:306:PRO:HD3	2.00	0.43
1:A:216[B]:GLU:OE1	1:A:216[B]:GLU:HA	2.20	0.42
1:B:356:VAL:O	1:B:356:VAL:HG23	2.19	0.42
1:A:23:LEU:O	1:A:23:LEU:HD23	2.20	0.41
1:A:150[A]:ARG:HB3	1:A:150[A]:ARG:NH1	2.35	0.41
1:A:132:GLU:H	1:A:132:GLU:CD	2.28	0.41
1:B:291:MET:HE2	1:B:320[A]:MET:HG3	2.02	0.41
1:B:150[B]:ARG:NH1	4:B:499:HOH:O	2.53	0.41
1:A:23:LEU:HD13	1:B:20:LEU:CD2	2.48	0.40
1:A:333:ASP:O	1:A:337:GLN:HG3	2.21	0.40
1:B:178:SER:HB2	1:B:185:VAL:HG23	2.03	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	354/385 (92%)	349 (99%)	5 (1%)	0	100	100
1	B	328/385 (85%)	323 (98%)	5 (2%)	0	100	100
All	All	682/770 (89%)	672 (98%)	10 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	302/325 (93%)	301 (100%)	1 (0%)	86	87
1	B	280/325 (86%)	274 (98%)	6 (2%)	47	41
All	All	582/650 (90%)	575 (99%)	7 (1%)	65	61

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	261	ARG
1	B	56	ASN
1	B	150[A]	ARG
1	B	150[B]	ARG
1	B	191	THR
1	B	261	ARG
1	B	301	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	15	GLN
1	A	86	GLN
1	A	100	ASN
1	A	251	GLN
1	B	86	GLN
1	B	100	ASN

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 ⓘ

Of 6 ligands modelled in this entry, 4 are monoatomic - leaving 2 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	$\# Z > 2$
3	ACT	B	363	-	3,3,3	0.87	0	3,3,3	1.35	0
3	ACT	A	363	-	3,3,3	0.84	0	3,3,3	1.57	1 (33%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
3	A	363	ACT	OXT-C-O	-2.07	114.35	122.03

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	363	ACT	4	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

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, 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	345/385 (89%)	1.23	75 (21%)	2 2	10, 31, 158, 259	13 (3%)
1	B	327/385 (84%)	0.89	66 (20%)	3 3	9, 25, 139, 217	7 (2%)
All	All	672/770 (87%)	1.07	141 (20%)	2 2	9, 28, 146, 259	20 (2%)

All (141) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	24	ILE	11.7
1	A	28	ALA	11.3
1	A	23	LEU	11.1
1	A	13	VAL	10.9
1	A	27	ARG	9.9
1	A	21	LEU	9.1
1	A	25	SER	8.7
1	A	18	ILE	8.5
1	A	20	LEU	8.3
1	A	17	ASN	7.8
1	A	16	LEU	7.2
1	B	58	ILE	6.9
1	B	86	GLN	6.8
1	A	66	PHE	6.7
1	A	26	LYS	6.7
1	A	9	LEU	6.5
1	B	31	VAL	6.4
1	A	37	ILE	6.3
1	B	6	LEU	6.3
1	B	79	PHE	6.2
1	A	39	GLY	5.9
1	B	30	LEU	5.9
1	B	55	LEU	5.9
1	B	85	LEU	5.9

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Mol	Chain	Res	Type	RSRZ
1	A	10	ARG	5.9
1	A	65	PRO	5.8
1	B	24	ILE	5.8
1	B	61	ALA	5.7
1	A	19	ASP	5.6
1	B	65	PRO	5.4
1	B	81	ALA	5.4
1	B	60	ALA	5.3
1	B	18	ILE	5.3
1	B	28	ALA	5.3
1	B	25	SER	5.2
1	B	59	LEU	5.1
1	B	356	VAL	5.1
1	B	23	LEU	5.1
1	A	357	PRO	5.1
1	B	29	ASN	5.0
1	B	82	GLY	4.9
1	A	85	LEU	4.9
1	B	57	THR	4.8
1	A	22	GLU	4.8
1	B	32	GLN	4.8
1	B	75	PHE	4.7
1	B	54	MET	4.7
1	B	71	VAL	4.6
1	A	71	VAL	4.6
1	B	83	LEU	4.6
1	B	74	LEU	4.5
1	B	66	PHE	4.5
1	B	78	ILE	4.5
1	A	74	LEU	4.4
1	B	9	LEU	4.3
1	A	14	ASP	4.3
1	A	64	GLY	4.3
1	A	46	PHE	4.2
1	A	61	ALA	4.2
1	A	49	LEU	4.2
1	B	7	GLU	4.1
1	A	11	THR	4.1
1	A	34	ILE	4.0
1	B	21	LEU	4.0
1	B	70	THR	3.9
1	A	15	GLN	3.9

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Mol	Chain	Res	Type	RSRZ
1	B	68	ASP	3.9
1	B	13	VAL	3.7
1	A	67[A]	GLU	3.6
1	A	30	LEU	3.6
1	A	36	LYS	3.5
1	B	76	LYS	3.5
1	B	63	GLU	3.5
1	A	60	ALA	3.5
1	A	355	LEU	3.5
1	B	10	ARG	3.5
1	A	12	GLN	3.4
1	B	11	THR	3.4
1	A	356	VAL	3.4
1	A	62	ASN	3.4
1	A	70	THR	3.4
1	B	27	ARG	3.4
1	A	38	LYS	3.4
1	B	69	SER	3.4
1	B	80	LYS	3.3
1	B	16	LEU	3.3
1	B	20	LEU	3.3
1	B	355	LEU	3.3
1	A	31	VAL	3.2
1	B	53	GLU	3.2
1	A	82	GLY	3.2
1	A	8	GLU	3.2
1	B	62	ASN	3.1
1	B	19	ASP	3.1
1	A	33	GLU	3.1
1	B	56	ASN	3.1
1	B	72	GLN	3.1
1	A	63	GLU	3.1
1	A	45	ARG	3.1
1	A	32	GLN	3.0
1	A	102	LYS	3.0
1	A	78	ILE	3.0
1	A	86	GLN	3.0
1	A	68	ASP	3.0
1	B	26	LYS	3.0
1	A	59	LEU	2.9
1	A	69	SER	2.9
1	A	354	ASN	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	331	LEU	2.8
1	B	8	GLU	2.8
1	B	73	LYS	2.7
1	B	15	GLN	2.7
1	A	75	PHE	2.7
1	B	67	GLU	2.7
1	B	345	GLU	2.7
1	B	64	GLY	2.6
1	B	91	SER	2.6
1	A	29	ASN	2.6
1	A	89	ASP	2.5
1	B	14	ASP	2.5
1	A	58	ILE	2.5
1	A	73	LYS	2.5
1	A	35	GLY	2.4
1	A	84	GLU	2.4
1	B	84	GLU	2.4
1	A	83	LEU	2.3
1	B	93	ALA	2.3
1	A	196	GLU	2.3
1	A	197	VAL	2.3
1	B	351	LEU	2.3
1	A	48	PRO	2.2
1	A	56	ASN	2.2
1	A	81	ALA	2.2
1	B	77	GLU	2.2
1	B	12	GLN	2.2
1	A	47	ASP	2.1
1	A	352	ALA	2.1
1	A	131	TYR	2.1
1	A	92	LYS	2.1
1	B	99	LYS	2.1
1	A	57	THR	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](#)

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, 95th 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 ²)	Q<0.9
3	ACT	B	363	4/4	0.80	0.15	36,38,38,39	0
3	ACT	A	363	4/4	0.84	0.15	46,46,46,47	0
2	MN	A	362[A]	1/1	0.99	0.06	18,18,18,18	1
2	MN	A	362[B]	1/1	0.99	0.06	29,29,29,29	1
2	MN	B	362[A]	1/1	1.00	0.04	18,18,18,18	1
2	MN	B	362[B]	1/1	1.00	0.04	22,22,22,22	1

6.5 Other polymers [i](#)

There are no such residues in this entry.