



Full wwPDB X-ray Structure Validation Report ⓘ

May 29, 2020 – 06:48 am BST

PDB ID : 5F7P
Title : Rok Repressor Lmo0178 from *Listeria monocytogenes*
Authors : Light, S.H.; Anderson, W.F.; Center for Structural Genomics of Infectious Diseases (CSGID)
Deposited on : 2015-12-08
Resolution : 2.84 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.11
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)	:	2.11

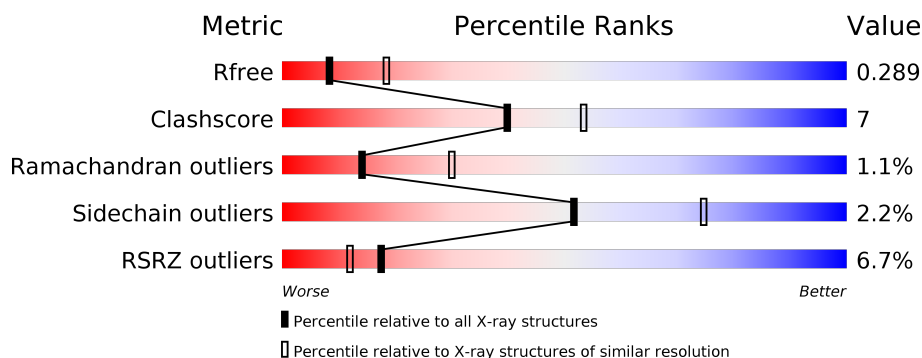
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.84 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	1031 (2.86-2.82)
Clashscore	141614	1078 (2.86-2.82)
Ramachandran outliers	138981	1050 (2.86-2.82)
Sidechain outliers	138945	1051 (2.86-2.82)
RSRZ outliers	127900	1019 (2.86-2.82)

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 ≥ 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	407	<div> <div>3%</div> <div> <div></div> <div>70%</div> <div>18%</div> <div>•</div> <div>12%</div> </div> </div>
1	E	407	<div> <div>7%</div> <div> <div></div> <div>60%</div> <div>13%</div> <div>27%</div> </div> </div>

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 5122 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 Lmo0178 protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	E	299	Total	C	N	O	S	0	0	0
			2312	1480	384	436	12			
1	A	359	Total	C	N	O	S	0	0	0
			2805	1785	472	533	15			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	-2	SER	-	expression tag	UNP Q8YAF1
E	-1	ASN	-	expression tag	UNP Q8YAF1
E	0	ALA	-	expression tag	UNP Q8YAF1
A	-2	SER	-	expression tag	UNP Q8YAF1
A	-1	ASN	-	expression tag	UNP Q8YAF1
A	0	ALA	-	expression tag	UNP Q8YAF1

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Zn	0	0
			1	1		
2	E	1	Total	Zn	0	0
			1	1		

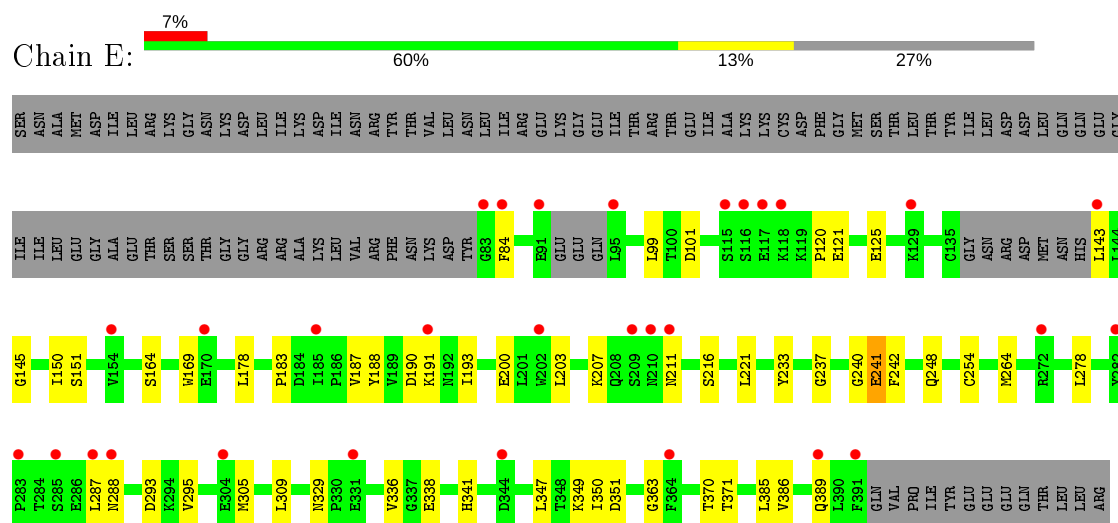
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	E	2	Total	O	0	0
			2	2		
3	A	1	Total	O	0	0
			1	1		

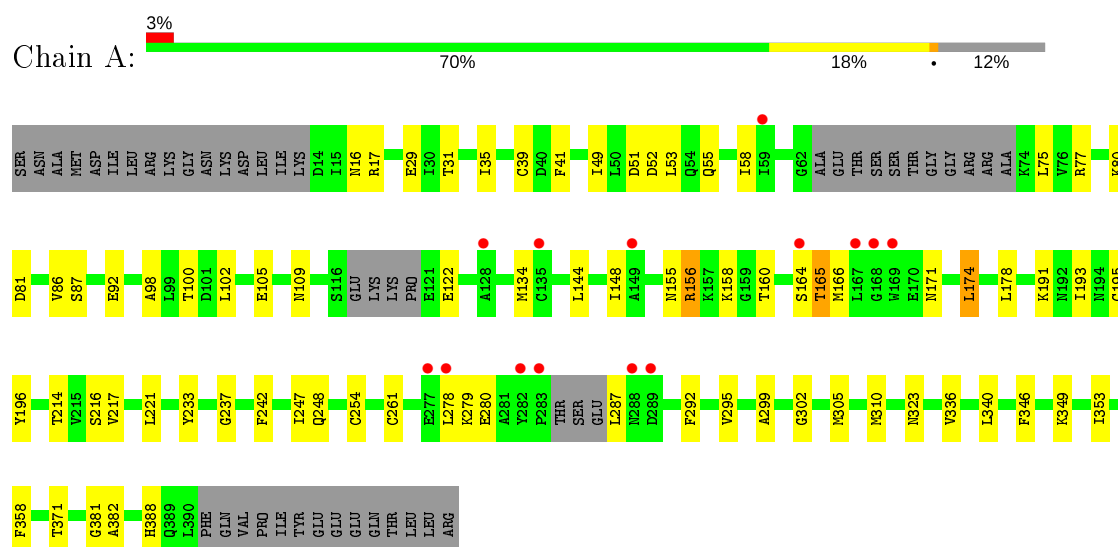
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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: Lmo0178 protein



- Molecule 1: Lmo0178 protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	68.97Å 81.78Å 141.09Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	70.76 – 2.84 29.32 – 2.84	Depositor EDS
% Data completeness (in resolution range)	97.1 (70.76-2.84) 97.3 (29.32-2.84)	Depositor EDS
R_{merge}	0.03	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.39 (at 2.85Å)	Xtriage
Refinement program	REFMAC 5.8.0135	Depositor
R, R_{free}	0.223 , 0.293 0.224 , 0.289	Depositor DCC
R_{free} test set	955 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	88.2	Xtriage
Anisotropy	0.039	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 59.2	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	0.93	EDS
Total number of atoms	5122	wwPDB-VP
Average B, all atoms (Å ²)	104.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.21% 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: ZN

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	0.73	0/2853	0.89	0/3845
1	E	0.63	0/2357	0.76	0/3181
All	All	0.68	0/5210	0.83	0/7026

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	2805	0	2776	45	0
1	E	2312	0	2286	31	0
2	A	1	0	0	0	0
2	E	1	0	0	0	0
3	A	1	0	0	0	0
3	E	2	0	0	0	0
All	All	5122	0	5062	74	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 (74) 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:247:ILE:O	1:A:247:ILE:HG13	1.88	0.71
1:A:39:CYS:SG	1:A:41:PHE:CZ	2.86	0.69
1:A:254:CYS:HB2	1:A:261:CYS:SG	2.36	0.66
1:E:248:GLN:HE22	1:A:248:GLN:HE22	1.45	0.64
1:A:148:ILE:HG21	1:A:174:LEU:HD21	1.79	0.63
1:E:336:VAL:HG12	1:E:371:THR:HB	1.79	0.62
1:A:295:VAL:HG12	1:A:310:MET:CE	2.30	0.62
1:E:200:GLU:OE2	1:E:371:THR:OG1	2.12	0.61
1:E:287:LEU:HD21	1:E:295:VAL:HG22	1.82	0.61
1:A:29:GLU:HG2	1:A:77:ARG:HB3	1.86	0.58
1:E:347:LEU:HD11	1:E:370:THR:HG21	1.86	0.57
1:A:164:SER:O	1:A:165:THR:C	2.45	0.55
1:E:193:ILE:HD13	1:E:216:SER:HB3	1.88	0.55
1:E:233:TYR:HB3	1:E:237:GLY:HA2	1.88	0.54
1:A:156:ARG:HD3	1:A:233:TYR:CE1	2.44	0.53
1:E:241:GLU:HG3	1:A:358:PHE:CD2	2.44	0.52
1:E:240:GLY:O	1:E:242:PHE:N	2.36	0.52
1:E:278:LEU:HD12	1:E:309:LEU:HD21	1.92	0.52
1:A:292:PHE:CD1	1:A:340:LEU:HD22	2.44	0.52
1:A:287:LEU:HD23	1:A:305:MET:CE	2.40	0.52
1:E:188:TYR:CE1	1:E:389:GLN:HG2	2.45	0.51
1:A:80:LYS:HE2	1:A:81:ASP:OD2	2.11	0.51
1:A:53:LEU:HD22	1:A:58:ILE:HG21	1.92	0.50
1:A:86:VAL:HG13	1:A:98:ALA:O	2.11	0.49
1:A:49:ILE:O	1:A:52:ASP:OD1	2.31	0.48
1:A:102:LEU:HD11	1:A:144:LEU:HD11	1.96	0.48
1:A:233:TYR:HB3	1:A:237:GLY:HA2	1.96	0.48
1:A:292:PHE:CG	1:A:340:LEU:HD22	2.49	0.48
1:A:254:CYS:CB	1:A:261:CYS:SG	2.98	0.47
1:A:92:GLU:CD	1:A:166:MET:SD	2.93	0.47
1:E:178:LEU:HD12	1:E:187:VAL:HG11	1.96	0.47
1:A:155:ASN:O	1:A:158:LYS:N	2.48	0.47
1:E:164:SER:HB3	1:E:169:TRP:HB2	1.96	0.47
1:E:254:CYS:HA	1:E:264:MET:SD	2.55	0.47
1:A:148:ILE:HD11	1:A:178:LEU:HD22	1.96	0.46
1:A:122:GLU:HG2	1:A:122:GLU:O	2.16	0.46
1:E:288:ASN:OD1	1:E:288:ASN:N	2.48	0.46
1:A:195:CYS:O	1:A:381:GLY:HA3	2.17	0.45
1:E:338:GLU:HG3	1:E:341:HIS:HD2	1.81	0.45
1:A:100:THR:HA	1:A:105:GLU:O	2.16	0.45
1:A:31:THR:HG22	1:A:75:LEU:HD12	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:84:PHE:CE1	1:E:143:LEU:N	2.86	0.44
1:E:121:GLU:O	1:E:125:GLU:HG2	2.17	0.44
1:E:188:TYR:CZ	1:E:389:GLN:HG2	2.53	0.44
1:A:31:THR:O	1:A:35:ILE:HG13	2.18	0.43
1:A:98:ALA:HA	1:A:134:MET:HE3	2.00	0.43
1:E:101:ASP:N	1:E:101:ASP:OD1	2.49	0.43
1:E:287:LEU:HD23	1:E:305:MET:CE	2.48	0.43
1:E:190:ASP:HB2	1:E:385:LEU:HD21	2.01	0.43
1:A:16:ASN:O	1:A:17:ARG:C	2.57	0.43
1:A:51:ASP:O	1:A:55:GLN:HG2	2.18	0.43
1:E:329:ASN:CG	1:E:329:ASN:O	2.57	0.43
1:E:150:ILE:CG1	1:E:151:SER:N	2.82	0.42
1:E:190:ASP:OD1	1:E:191:LYS:N	2.48	0.42
1:A:165:THR:OG1	1:A:166:MET:N	2.53	0.42
1:A:242:PHE:HZ	1:A:323:ASN:HB2	1.85	0.42
1:A:336:VAL:HG22	1:A:371:THR:HB	2.02	0.42
1:A:39:CYS:SG	1:A:41:PHE:CE2	3.12	0.41
1:A:193:ILE:HD11	1:A:214:THR:HG22	2.01	0.41
1:A:299:ALA:O	1:A:302:GLY:N	2.33	0.41
1:E:203:LEU:O	1:E:207:LYS:CE	2.68	0.41
1:E:264:MET:HE2	1:E:264:MET:HA	2.01	0.41
1:E:371:THR:HG23	1:E:371:THR:O	2.20	0.41
1:A:196:TYR:O	1:A:196:TYR:CD1	2.73	0.41
1:E:349:LYS:O	1:E:350:ILE:C	2.57	0.41
1:E:145:GLY:HA3	1:E:386:VAL:HG11	2.01	0.41
1:A:53:LEU:HD22	1:A:58:ILE:CG2	2.50	0.41
1:A:109:ASN:HA	1:A:134:MET:HE3	2.02	0.41
1:A:87:SER:HB2	1:A:382:ALA:HB3	2.03	0.40
1:A:191:LYS:HE3	1:A:193:ILE:HG23	2.04	0.40
1:A:109:ASN:HA	1:A:134:MET:CE	2.51	0.40
1:E:84:PHE:CD2	1:E:99:LEU:HD21	2.57	0.40
1:A:278:LEU:O	1:A:279:LYS:C	2.58	0.40
1:A:346:PHE:HB3	1:A:349:LYS:HG2	2.03	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	351/407 (86%)	318 (91%)	30 (8%)	3 (1%)	17	34
1	E	293/407 (72%)	263 (90%)	26 (9%)	4 (1%)	11	24
All	All	644/814 (79%)	581 (90%)	56 (9%)	7 (1%)	14	30

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	241	GLU
1	A	171	ASN
1	E	183	PRO
1	E	363	GLY
1	A	156	ARG
1	A	165	THR
1	E	120	PRO

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	298/340 (88%)	290 (97%)	8 (3%)	44	69
1	E	244/340 (72%)	240 (98%)	4 (2%)	62	81
All	All	542/680 (80%)	530 (98%)	12 (2%)	52	75

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

Mol	Chain	Res	Type
1	E	211	ASN
1	E	221	LEU
1	E	293	ASP
1	E	351	ASP
1	A	160	THR
1	A	174	LEU
1	A	216	SER
1	A	217	VAL
1	A	221	LEU
1	A	280	GLU
1	A	353	ILE
1	A	388	HIS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	E	130	ASN
1	E	230	GLN
1	E	248	GLN
1	E	258	GLN
1	A	103	ASN
1	A	142	HIS
1	A	208	GLN
1	A	356	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 carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	359/407 (88%)	0.09	14 (3%) 39 31	48, 93, 138, 160	0
1	E	299/407 (73%)	0.33	30 (10%) 7 4	51, 108, 171, 194	0
All	All	658/814 (80%)	0.20	44 (6%) 17 12	48, 100, 154, 194	0

All (44) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	202	TRP	6.9
1	E	116	SER	5.7
1	A	278	LEU	5.5
1	E	143	LEU	5.4
1	E	83	GLY	5.2
1	E	391	PHE	5.2
1	A	282	TYR	4.6
1	E	282	TYR	4.5
1	E	287	LEU	3.9
1	E	84	PHE	3.9
1	E	331	GLU	3.8
1	A	167	LEU	3.7
1	E	210	ASN	3.5
1	E	115	SER	3.4
1	E	117	GLU	3.3
1	E	211	ASN	3.2
1	E	209	SER	3.2
1	A	164	SER	3.1
1	A	277	GLU	3.0
1	E	389	GLN	2.9
1	E	288	ASN	2.8
1	E	185	ILE	2.8
1	E	364	PHE	2.8
1	A	135	CYS	2.8

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Mol	Chain	Res	Type	RSRZ
1	E	272	ARG	2.7
1	A	289	ASP	2.7
1	A	168	GLY	2.7
1	E	344	ASP	2.7
1	E	191	LYS	2.6
1	E	170	GLU	2.6
1	E	283	PRO	2.6
1	A	59	ILE	2.5
1	E	129	LYS	2.5
1	E	118	LYS	2.5
1	E	304	GLU	2.4
1	A	288	ASN	2.4
1	A	149	ALA	2.4
1	A	128	ALA	2.2
1	A	169	TRP	2.2
1	A	283	PRO	2.2
1	E	91	GLU	2.2
1	E	95	LEU	2.1
1	E	285	SER	2.0
1	E	154	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 carbohydrates 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(\AA^2)	Q<0.9
2	ZN	A	501	1/1	0.99	0.08	69,69,69,69	0
2	ZN	E	501	1/1	0.99	0.10	86,86,86,86	0

6.5 Other polymers [i](#)

There are no such residues in this entry.