



Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 03:08 AM GMT

PDB ID : 3R3L
Title : Structure of NP protein from Lassa AV strain
Authors : Perbandt, M.; Brunotte, L.; Gunther, S.; Betzel, C.
Deposited on : 2011-03-16
Resolution : 2.45 Å(reported)

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

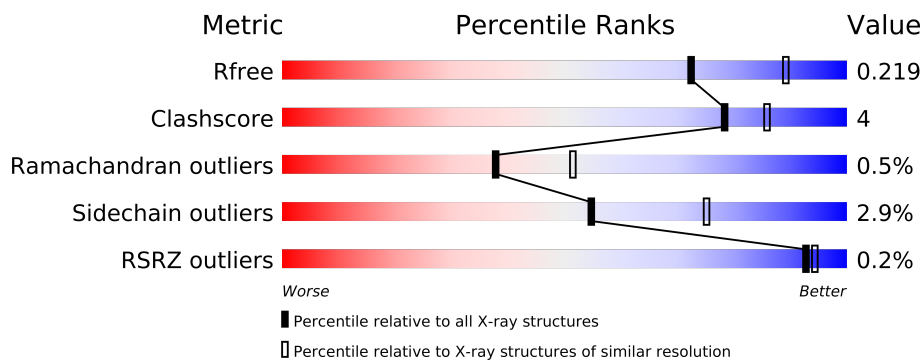
The following versions of software and data (see [references](#)) were used in the production of this report:

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 2.45 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	2989 (2.48-2.40)
Clashscore	79885	3698 (2.48-2.40)
Ramachandran outliers	78287	3639 (2.48-2.40)
Sidechain outliers	78261	3640 (2.48-2.40)
RSRZ outliers	66119	2993 (2.48-2.40)

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. 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	582	
1	B	582	
1	C	582	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 12140 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 Nucleoprotein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	513	Total	C	N	O	S	0	0	0
			4009	2522	698	763	26			
1	B	514	Total	C	N	O	S	0	0	0
			4018	2528	700	764	26			
1	C	510	Total	C	N	O	S	0	0	0
			3986	2508	694	758	26			

There are 39 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	570	ASP	-	EXPRESSION TAG	UNP Q9DQX7
A	571	TYR	-	EXPRESSION TAG	UNP Q9DQX7
A	572	LYS	-	EXPRESSION TAG	UNP Q9DQX7
A	573	ASP	-	EXPRESSION TAG	UNP Q9DQX7
A	574	HIS	-	EXPRESSION TAG	UNP Q9DQX7
A	575	ASP	-	EXPRESSION TAG	UNP Q9DQX7
A	576	GLY	-	EXPRESSION TAG	UNP Q9DQX7
A	577	HIS	-	EXPRESSION TAG	UNP Q9DQX7
A	578	HIS	-	EXPRESSION TAG	UNP Q9DQX7
A	579	HIS	-	EXPRESSION TAG	UNP Q9DQX7
A	580	HIS	-	EXPRESSION TAG	UNP Q9DQX7
A	581	HIS	-	EXPRESSION TAG	UNP Q9DQX7
A	582	HIS	-	EXPRESSION TAG	UNP Q9DQX7
B	570	ASP	-	EXPRESSION TAG	UNP Q9DQX7
B	571	TYR	-	EXPRESSION TAG	UNP Q9DQX7
B	572	LYS	-	EXPRESSION TAG	UNP Q9DQX7
B	573	ASP	-	EXPRESSION TAG	UNP Q9DQX7
B	574	HIS	-	EXPRESSION TAG	UNP Q9DQX7
B	575	ASP	-	EXPRESSION TAG	UNP Q9DQX7
B	576	GLY	-	EXPRESSION TAG	UNP Q9DQX7
B	577	HIS	-	EXPRESSION TAG	UNP Q9DQX7
B	578	HIS	-	EXPRESSION TAG	UNP Q9DQX7
B	579	HIS	-	EXPRESSION TAG	UNP Q9DQX7

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Chain	Residue	Modelled	Actual	Comment	Reference
B	580	HIS	-	EXPRESSION TAG	UNP Q9DQX7
B	581	HIS	-	EXPRESSION TAG	UNP Q9DQX7
B	582	HIS	-	EXPRESSION TAG	UNP Q9DQX7
C	570	ASP	-	EXPRESSION TAG	UNP Q9DQX7
C	571	TYR	-	EXPRESSION TAG	UNP Q9DQX7
C	572	LYS	-	EXPRESSION TAG	UNP Q9DQX7
C	573	ASP	-	EXPRESSION TAG	UNP Q9DQX7
C	574	HIS	-	EXPRESSION TAG	UNP Q9DQX7
C	575	ASP	-	EXPRESSION TAG	UNP Q9DQX7
C	576	GLY	-	EXPRESSION TAG	UNP Q9DQX7
C	577	HIS	-	EXPRESSION TAG	UNP Q9DQX7
C	578	HIS	-	EXPRESSION TAG	UNP Q9DQX7
C	579	HIS	-	EXPRESSION TAG	UNP Q9DQX7
C	580	HIS	-	EXPRESSION TAG	UNP Q9DQX7
C	581	HIS	-	EXPRESSION TAG	UNP Q9DQX7
C	582	HIS	-	EXPRESSION TAG	UNP Q9DQX7

- Molecule 2 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

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

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

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

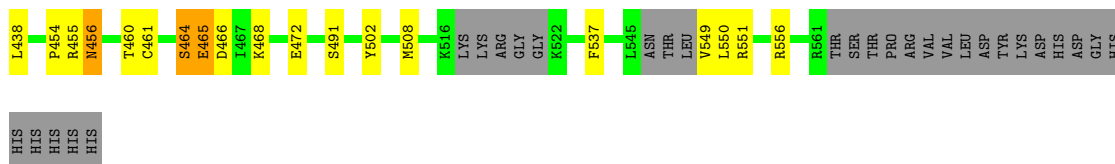
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	41	Total O 41 41	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	48	Total 48	O 48	0	0
4	C	30	Total 30	O 30	0	0



4 Data and refinement statistics

Property	Value	Source
Space group	P 3	Depositor
Cell constants a, b, c, α , β , γ	176.32Å 176.32Å 56.58Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	30.00 – 2.45 88.16 – 2.44	Depositor EDS
% Data completeness (in resolution range)	51.4 (30.00-2.45) 51.4 (88.16-2.44)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.28 (at 2.45Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.184 , 0.219 0.182 , 0.219	Depositor DCC
R_{free} test set	1873 reflections (5.25%)	DCC
Wilson B-factor (Å ²)	33.2	Xtriage
Anisotropy	0.331	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , -16.9	EDS
Estimated twinning fraction	0.332 for H, K, L 0.163 for -H, H+K, -L 0.163 for -H, -K, L 0.342 for -H-K, K, -L 0.315 for -h,-k,l 0.499 for h,-h-k,-l 0.354 for -k,-h,-l	Xtriage
Reported twinning fraction	0.332 for H, K, L 0.163 for -H, H+K, -L 0.163 for -H, -K, L 0.342 for -H-K, K, -L	Depositor
L-test for twinning	$\langle L \rangle = 0.37$, $\langle L^2 \rangle = 0.19$	Xtriage
Outliers	0 of 37572 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	12140	wwPDB-VP
Average B, all atoms (Å ²)	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.72% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities 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: ZN, MN

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	0.32	0/4066	0.51	0/5489
1	B	0.31	0/4075	0.50	0/5500
1	C	0.34	0/4042	0.52	0/5454
All	All	0.32	0/12183	0.51	0/16443

There are no bond length outliers.

There are no bond angle outliers.

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	4009	0	4089	20	0
1	B	4018	0	4102	29	0
1	C	3986	0	4064	43	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
2	C	1	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	41	0	0	0	0
4	B	48	0	0	0	0
4	C	30	0	0	0	0
All	All	12140	0	12255	88	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 4.

All (88) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:216:THR:HA	1:A:234:THR:HG21	1.73	0.70
1:C:455:ARG:O	1:C:456:ASN:HB2	1.94	0.67
1:B:196:ASN:O	1:B:200:GLN:HG2	1.99	0.63
1:B:420:LEU:HD13	1:B:435:ILE:HD12	1.85	0.58
1:C:81:LYS:HA	1:C:322:SER:H	1.68	0.58
1:A:89:LEU:HD12	1:A:336:VAL:HG22	1.87	0.56
1:C:416:GLU:HG3	1:C:438:LEU:O	2.05	0.56
1:C:47:VAL:HG22	1:C:70:LEU:HD21	1.87	0.55
1:C:373:LEU:O	1:C:377:MET:HG2	2.07	0.55
1:B:187:LEU:HB3	1:B:228:ILE:HD11	1.88	0.55
1:A:49:ASN:O	1:A:53:LEU:HG	2.07	0.54
1:A:17:ARG:HG2	1:A:263:ASN:O	2.08	0.54
1:A:221:ARG:HE	1:B:207:LEU:HD21	1.73	0.53
1:B:233:ASP:HB3	1:B:236:LYS:HG3	1.90	0.53
1:C:464:SER:O	1:C:466:ASP:N	2.42	0.53
1:C:122:SER:HA	1:C:551:ARG:HD3	1.90	0.53
1:B:223:THR:HG23	1:B:230:ASN:OD1	2.08	0.53
1:B:382:PRO:HB3	1:B:454:PRO:HB3	1.91	0.53
1:C:388:MET:HG3	1:C:402:LEU:HG	1.92	0.52
1:B:180:PRO:HB3	1:B:254:ALA:HA	1.91	0.52
1:C:41:GLY:O	1:C:77:LEU:HD11	2.10	0.51
1:C:196:ASN:O	1:C:200:GLN:HG2	2.11	0.51
1:B:420:LEU:O	1:B:424:LYS:HG2	2.10	0.51
1:C:382:PRO:HB3	1:C:454:PRO:CB	2.41	0.50
1:C:17:ARG:HG2	1:C:265:LEU:HG	1.94	0.50
1:B:484:ILE:HG22	1:B:486:LEU:HG	1.94	0.49
1:A:196:ASN:O	1:A:200:GLN:HG2	2.11	0.49
1:C:27:ILE:HD11	1:C:267:THR:HG22	1.95	0.49
1:C:468:LYS:O	1:C:472:GLU:HB2	2.11	0.49
1:C:14:GLN:HA	1:C:17:ARG:HD2	1.95	0.48
1:C:9:SER:O	1:C:302:PRO:HG2	2.13	0.48
1:C:234:THR:HA	1:C:248:LEU:HD13	1.95	0.48

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:204:ASP:O	1:B:208:ILE:HG12	2.14	0.48
1:A:409:CYS:HA	1:A:551:ARG:O	2.14	0.48
1:C:396:ASP:HA	1:C:415:ARG:HH12	1.79	0.47
1:A:317:TRP:HB2	1:A:321:ALA:HB2	1.96	0.47
1:C:71:ASN:HB3	1:C:324:THR:HB	1.96	0.47
1:C:239:LEU:C	1:C:241:ILE:H	2.18	0.47
1:B:440:ALA:O	1:B:561:ARG:HG2	2.14	0.47
1:A:78:VAL:HG11	1:A:179:MET:HE3	1.96	0.47
1:C:502:TYR:CE1	1:C:550:LEU:HD13	2.49	0.47
1:B:491:SER:HB2	1:B:537:PHE:CE1	2.50	0.47
1:C:38:LEU:HD22	1:C:42:LEU:HD22	1.95	0.47
1:B:216:THR:HA	1:B:234:THR:HG21	1.96	0.47
1:A:365:LEU:HD23	1:A:369:GLN:HB3	1.97	0.47
1:B:367:TYR:O	1:B:371:MET:HG3	2.15	0.47
1:B:489:ILE:HG13	1:B:489:ILE:H	1.54	0.47
1:C:387:TRP:HA	1:C:460:THR:O	2.16	0.46
1:B:36:GLN:HE21	1:B:195:LEU:CD2	2.28	0.46
1:A:204:ASP:O	1:A:208:ILE:HG12	2.15	0.46
1:B:220:ASP:O	1:B:223:THR:HB	2.16	0.46
1:C:286:LYS:HA	1:C:291:MET:HB2	1.98	0.45
1:B:317:TRP:HB2	1:B:321:ALA:HB2	1.99	0.45
1:C:180:PRO:HB3	1:C:254:ALA:HA	1.98	0.45
1:A:180:PRO:HG2	1:A:320:ILE:HG22	1.99	0.44
1:C:408:GLY:O	1:C:550:LEU:HA	2.17	0.44
1:C:217:SER:HA	1:C:220:ASP:HB2	2.00	0.43
1:C:25:SER:HB2	1:C:26:ASN:H	1.69	0.43
1:A:207:LEU:HD11	1:B:221:ARG:HB2	2.00	0.43
1:C:420:LEU:HD11	1:C:424:LYS:HE2	2.00	0.43
1:A:468:LYS:O	1:A:472:GLU:HB2	2.17	0.43
1:A:207:LEU:HD21	1:B:221:ARG:HE	1.84	0.43
1:B:20:LEU:HD22	1:B:278:ILE:HG23	1.99	0.43
1:C:236:LYS:N	1:C:236:LYS:HD3	2.34	0.43
1:B:9:SER:O	1:B:302:PRO:HG2	2.19	0.43
1:C:223:THR:HA	1:C:226:HIS:O	2.18	0.43
1:C:382:PRO:HB3	1:C:454:PRO:HB2	2.01	0.42
1:C:491:SER:HB2	1:C:537:PHE:CE1	2.54	0.42
1:B:54:MET:O	1:B:59:ARG:NH2	2.51	0.42
1:C:178:THR:HG22	1:C:243:GLY:HA2	2.01	0.42
1:C:380:LEU:HD11	1:C:411:ILE:HD11	2.02	0.42
1:C:455:ARG:O	1:C:456:ASN:CB	2.66	0.42
1:C:316:GLY:O	1:C:318:PRO:HD3	2.19	0.42
1:A:426:ASP:O	1:A:430:SER:N	2.50	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:296:THR:HA	1:C:297:PRO:HD3	1.89	0.42
1:B:44:PHE:CG	1:B:189:LYS:HG2	2.55	0.42
1:C:437:ASP:HB3	1:C:508:MET:SD	2.61	0.41
1:A:309:LYS:HE3	1:A:323:ARG:HH12	1.86	0.41
1:C:39:LEU:HD11	1:C:195:LEU:HD13	2.01	0.41
1:B:420:LEU:HG	1:B:424:LYS:HE2	2.03	0.41
1:B:209:TYR:HA	1:B:212:LYS:O	2.21	0.41
1:A:210:THR:HG22	1:A:264:MET:CE	2.50	0.41
1:B:19:GLU:HG3	1:B:281:SER:HB3	2.03	0.40
1:C:111:SER:O	1:C:115:ARG:HG3	2.22	0.40
1:A:380:LEU:HD11	1:A:411:ILE:HD11	2.03	0.40
1:A:225:SER:HB3	1:B:200:GLN:HB3	2.04	0.40
1:C:38:LEU:HD21	1:C:74:VAL:HG13	2.04	0.40
1:C:401:ALA:HA	1:C:411:ILE:O	2.21	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	505/582 (87%)	486 (96%)	19 (4%)	0	100	100
1	B	506/582 (87%)	490 (97%)	16 (3%)	0	100	100
1	C	500/582 (86%)	465 (93%)	28 (6%)	7 (1%)	16	21
All	All	1511/1746 (86%)	1441 (95%)	63 (4%)	7 (0%)	38	51

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	91	VAL
1	C	464	SER
1	C	465	GLU
1	C	27	ILE
1	C	240	ASN

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Mol	Chain	Res	Type
1	C	456	ASN
1	C	92	GLY

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	450/505 (89%)	441 (98%)	9 (2%)	68	85
1	B	451/505 (89%)	439 (97%)	12 (3%)	57	77
1	C	447/505 (88%)	429 (96%)	18 (4%)	42	62
All	All	1348/1515 (89%)	1309 (97%)	39 (3%)	55	75

All (39) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	98	ASP
1	A	233	ASP
1	A	235	LYS
1	A	263	ASN
1	A	395	GLU
1	A	465	GLU
1	A	466	ASP
1	A	545	LEU
1	A	546	ASN
1	B	8	LYS
1	B	25	SER
1	B	83	THR
1	B	90	ARG
1	B	112	LYS
1	B	131	SER
1	B	158	ASP
1	B	234	THR
1	B	296	THR
1	B	370	LEU
1	B	464	SER
1	B	489	ILE

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Mol	Chain	Res	Type
1	C	8	LYS
1	C	25	SER
1	C	28	LYS
1	C	83	THR
1	C	107	GLU
1	C	112	LYS
1	C	235	LYS
1	C	236	LYS
1	C	264	MET
1	C	271	SER
1	C	287	ARG
1	C	323	ARG
1	C	327	VAL
1	C	367	TYR
1	C	461	CYS
1	C	465	GLU
1	C	549	VAL
1	C	556	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	75	ASN
1	A	501	GLN
1	B	36	GLN
1	B	263	ASN
1	C	263	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 8 ligands modelled in this entry, 8 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.

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, 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	513/582 (88%)	0.05	2 (0%) 90 91	20, 20, 20, 26	1 (0%)
1	B	514/582 (88%)	0.09	0 100 100	20, 20, 20, 20	0
1	C	510/582 (87%)	-0.06	1 (0%) 93 94	20, 20, 20, 30	0
All	All	1537/1746 (88%)	0.03	3 (0%) 93 94	20, 20, 20, 30	1 (0%)

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	556	ARG	2.4
1	C	27	ILE	2.2
1	A	547	THR	2.0

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, 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	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	MN	C	583	1/1	0.06	-	30,30,30,30	0
3	ZN	C	584	1/1	0.09	-	20,20,20,20	0
3	ZN	A	585	1/1	0.16	-	20,20,20,20	0
3	ZN	B	585	1/1	0.17	-	20,20,20,20	0
2	MN	B	584	1/1	0.07	-	30,30,30,30	0
2	MN	A	584	1/1	0.09	-	30,30,30,30	0
2	MN	A	583	1/1	0.06	-	30,30,30,30	0
2	MN	B	583	1/1	0.09	-	30,30,30,30	0

6.5 Other polymers

There are no such residues in this entry.