



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 12:22 PM 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 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  
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>  
with specific help available everywhere you see the ⓘ symbol.

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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 : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

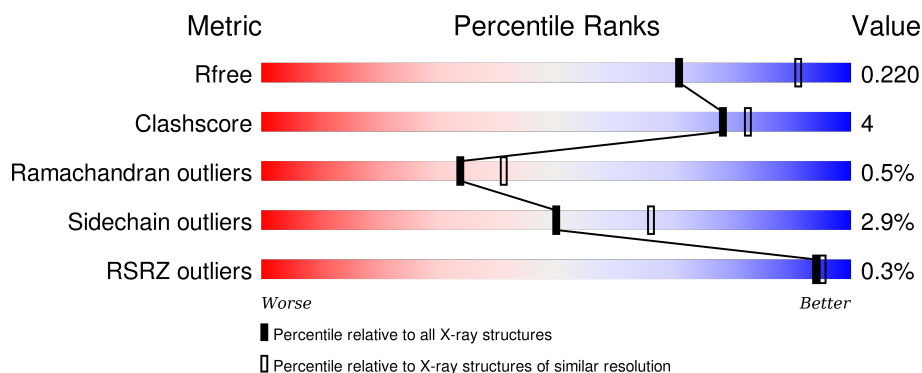
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.45 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))
$R_{free}$	91344	1003 (2.46-2.42)
Clashscore	102246	1071 (2.46-2.42)
Ramachandran outliers	100387	1065 (2.46-2.42)
Sidechain outliers	100360	1065 (2.46-2.42)
RSRZ outliers	91569	1005 (2.46-2.42)

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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. 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	582	 81% 7% 12%
1	B	582	 79% 9% 12%
1	C	582	 71% 15% • 12%

## 2 Entry composition

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

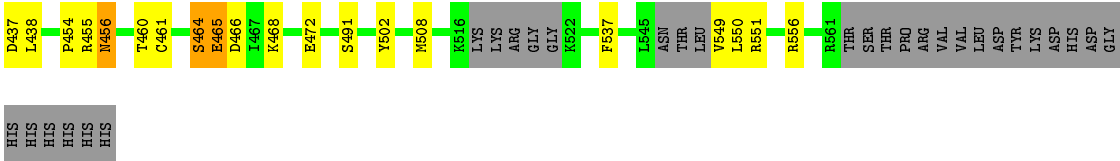
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	41	Total O 41 41	0	0

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





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 3	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	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
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.28 (at 2.45Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, $R_{free}$	0.184 , 0.219 0.182 , 0.220	Depositor DCC
$R_{free}$ test set	1862 reflections (5.26%)	DCC
Wilson B-factor (Å <sup>2</sup> )	33.2	Xtriage
Anisotropy	0.331	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , -16.1	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 <sup>2</sup>	$\langle  L  \rangle = 0.37$ , $\langle L^2 \rangle = 0.19$	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 <sup>2</sup> )	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.*

<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, 0.375 respectively for untwinned datasets, and 0.333, 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, 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 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	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
4	A	41	0	0	0	0

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

All (88) 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: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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:234:THR:HA	1:C:248:LEU:HD13	1.95	0.48
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
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 [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	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%)	14	13
All	All	1511/1746 (86%)	1441 (95%)	63 (4%)	7 (0%)	34	41

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	91	VAL

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Mol	Chain	Res	Type
1	C	464	SER
1	C	465	GLU
1	C	27	ILE
1	C	240	ASN
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%)	63	77
1	B	451/505 (89%)	439 (97%)	12 (3%)	52	69
1	C	447/505 (88%)	429 (96%)	18 (4%)	38	53
All	All	1348/1515 (89%)	1309 (97%)	39 (3%)	50	66

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

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Mol	Chain	Res	Type
1	B	234	THR
1	B	296	THR
1	B	370	LEU
1	B	464	SER
1	B	489	ILE
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 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 [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

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.

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 [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	513/582 (88%)	0.02	2 (0%) 93 94	20, 20, 20, 26	1 (0%)
1	B	514/582 (88%)	0.06	1 (0%) 95 96	20, 20, 20, 20	0
1	C	510/582 (87%)	-0.10	2 (0%) 93 94	20, 20, 20, 30	0
All	All	1537/1746 (88%)	-0.00	5 (0%) 94 95	20, 20, 20, 30	1 (0%)

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	556	ARG	2.6
1	C	27	ILE	2.1
1	A	547	THR	2.1
1	B	207	LEU	2.1
1	C	237	SER	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. 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, 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	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
3	ZN	A	585	1/1	1.00	0.16	0.71	20,20,20,20	0
3	ZN	B	585	1/1	0.98	0.17	-0.10	20,20,20,20	0
3	ZN	C	584	1/1	0.98	0.09	-1.21	20,20,20,20	0
2	MN	C	583	1/1	0.98	0.06	-2.57	30,30,30,30	0
2	MN	A	583	1/1	0.98	0.06	-2.99	30,30,30,30	0
2	MN	B	583	1/1	0.97	0.10	-3.72	30,30,30,30	0
2	MN	B	584	1/1	0.95	0.07	-4.97	30,30,30,30	0
2	MN	A	584	1/1	0.99	0.09	-7.04	30,30,30,30	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.