



# Full wwPDB X-ray Structure Validation Report ⓘ

Aug 21, 2020 – 03:27 PM BST

PDB ID : 6JNI  
Title : Crystal structure of the transcriptional regulator CadR from *P. putida* in complex with Zinc(II) and DNA  
Authors : Liu, X.C.; Gan, J.H.; Chen, H.  
Deposited on : 2019-03-16  
Resolution : 2.90 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

<https://www.wwpdb.org/validation/2017/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:

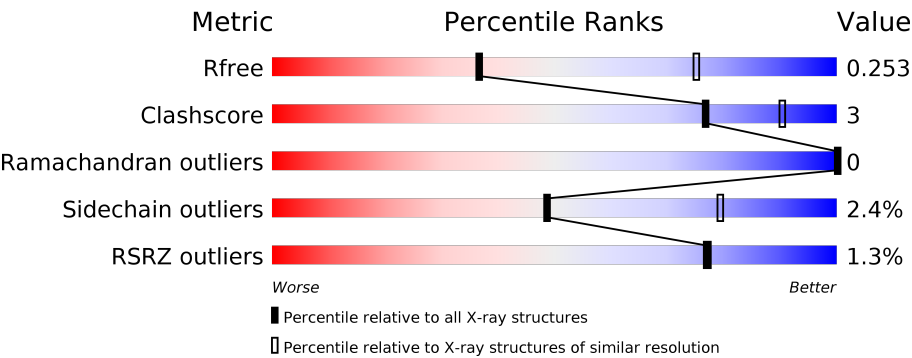
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.13.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.13.1

# 1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:  
*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.90 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.









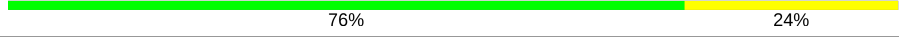

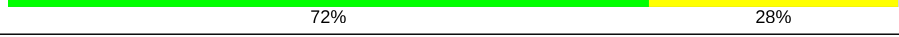

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	130704	1957 (2.90-2.90)
Clashscore	141614	2172 (2.90-2.90)
Ramachandran outliers	138981	2115 (2.90-2.90)
Sidechain outliers	138945	2117 (2.90-2.90)
RSRZ outliers	127900	1906 (2.90-2.90)

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 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	147	<div><div>2%</div><div><div></div><div>91%</div><div>7%</div><div></div></div><div></div></div>
1	B	147	<div><div></div><div><div></div><div>81%</div><div>5%</div><div>14%</div><div></div></div><div></div></div>
1	C	147	<div><div>%</div><div><div></div><div>84%</div><div></div><div>13%</div><div></div></div><div></div></div>
1	D	147	<div><div>%</div><div><div></div><div>91%</div><div>7%</div><div></div></div><div></div></div>
1	E	147	<div><div>3%</div><div><div></div><div>86%</div><div>6%</div><div>7%</div><div></div></div><div></div></div>
1	F	147	<div><div>%</div><div><div></div><div>82%</div><div>5%</div><div>13%</div><div></div></div><div></div></div>

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Mol	Chain	Length	Quality of chain
1	G	147	 84% 12%
1	H	147	 87% 6% 7%
2	I	25	 80% 16%
2	K	25	 80% 20%
2	M	25	 68% 20% 12%
2	O	25	 76% 24%
3	J	25	 76% 24%
3	L	25	 72% 20% 8%
3	N	25	 72% 28%
3	P	25	 60% 28% 8%

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 12608 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 CadR.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	145	Total	C	N	O	S	0	0	0
			1150	700	217	226	7			
1	B	126	Total	C	N	O	S	0	0	0
			1007	617	187	196	7			
1	C	128	Total	C	N	O	S	0	0	0
			1021	627	189	198	7			
1	D	145	Total	C	N	O	S	0	0	0
			1150	700	217	226	7			
1	E	137	Total	C	N	O	S	0	0	0
			1089	665	209	208	7			
1	F	128	Total	C	N	O	S	0	0	0
			1021	627	189	198	7			
1	G	129	Total	C	N	O	S	0	0	0
			1030	632	190	201	7			
1	H	137	Total	C	N	O	S	0	0	0
			1085	662	207	209	7			

- Molecule 2 is a DNA chain called DNA (25-MER).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	I	25	Total	C	N	O	P	0	0	0
			513	245	94	150	24			
2	K	25	Total	C	N	O	P	0	0	0
			497	235	92	146	24			
2	M	25	Total	C	N	O	P	0	0	0
			513	245	94	150	24			
2	O	25	Total	C	N	O	P	0	0	0
			513	245	94	150	24			

- Molecule 3 is a DNA chain called DNA (25-MER).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	J	25	Total	C	N	O	P	0	0	0
			506	242	94	146	24			
3	L	25	Total	C	N	O	P	0	0	0
			506	242	94	146	24			
3	N	25	Total	C	N	O	P	0	0	0
			506	242	94	146	24			
3	P	24	Total	C	N	O	P	0	0	0
			485	232	89	141	23			

- Molecule 4 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	G	2	Total	Zn	0	0
			2	2		
4	D	3	Total	Zn	0	0
			3	3		
4	E	3	Total	Zn	0	0
			3	3		
4	H	2	Total	Zn	0	0
			2	2		
4	B	1	Total	Zn	0	0
			1	1		
4	C	1	Total	Zn	0	0
			1	1		
4	A	3	Total	Zn	0	0
			3	3		
4	F	1	Total	Zn	0	0
			1	1		

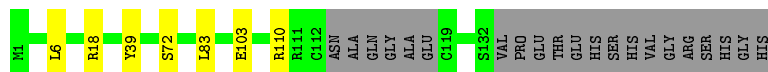
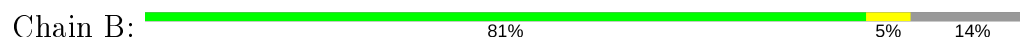
### 3 Residue-property plots [i](#)

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.

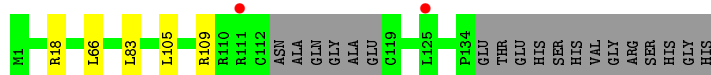
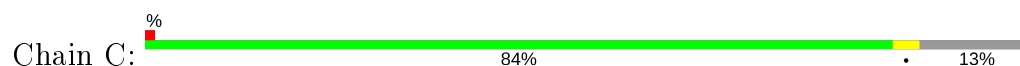
- Molecule 1: CadR



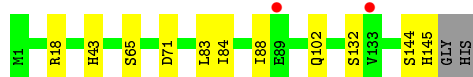
- Molecule 1: CadR



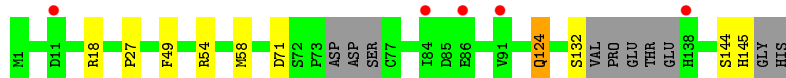
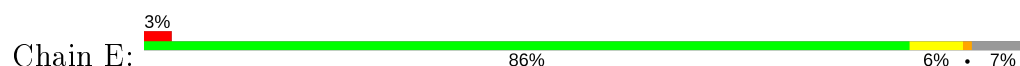
- Molecule 1: CadR



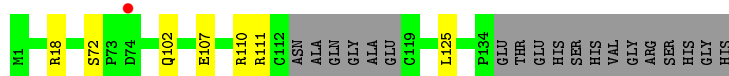
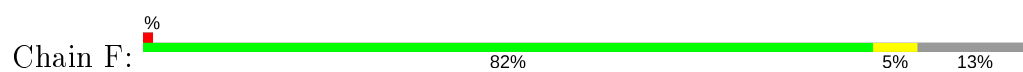
- Molecule 1: CadR



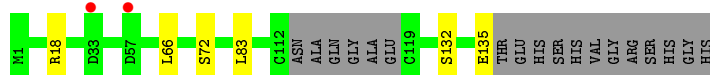
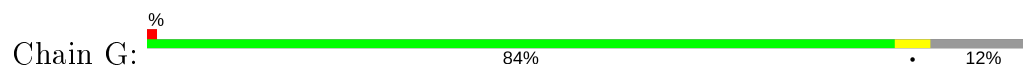
- Molecule 1: CadR



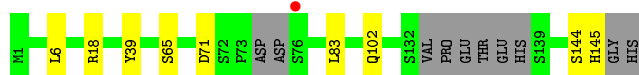
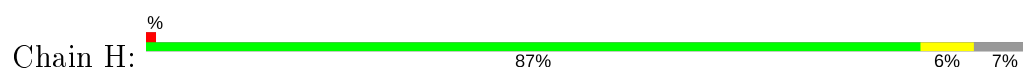
- Molecule 1: CadR



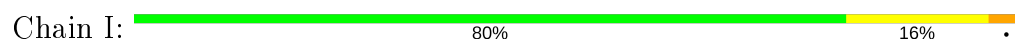
- Molecule 1: CadR



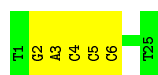
- Molecule 1: CadR



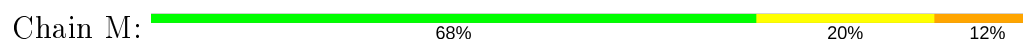
- Molecule 2: DNA (25-MER)



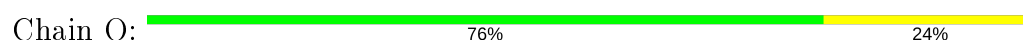
- Molecule 2: DNA (25-MER)



- Molecule 2: DNA (25-MER)



- Molecule 2: DNA (25-MER)



- Molecule 3: DNA (25-MER)

Chain J: 

76%

24%



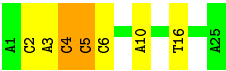
● Molecule 3: DNA (25-MER)

Chain L: 

72%

20%

8%



● Molecule 3: DNA (25-MER)

Chain N: 

72%

28%



● Molecule 3: DNA (25-MER)

Chain P: 

60%

28%

8%





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	71.09Å 71.33Å 98.80Å 81.03° 83.52° 73.85°	Depositor
Resolution (Å)	30.00 – 2.90 38.63 – 2.90	Depositor EDS
% Data completeness (in resolution range)	95.3 (30.00-2.90) 95.3 (38.63-2.90)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.12	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.45 (at 2.90Å)	Xtriage
Refinement program	REFMAC 5.8.0135	Depositor
R, $R_{free}$	0.212 , 0.254 0.214 , 0.253	Depositor DCC
$R_{free}$ test set	1939 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	55.9	Xtriage
Anisotropy	0.498	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.30 , 17.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.094 for -k,-h,-l	Xtriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	12608	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	62.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.43% 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, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

## 5 Model quality ⓘ

### 5.1 Standard geometry ⓘ

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	0.62	0/1164	0.83	0/1573
1	B	0.68	0/1016	0.86	0/1370
1	C	0.69	0/1031	0.83	0/1392
1	D	0.61	0/1164	0.84	0/1573
1	E	0.59	0/1100	0.81	0/1481
1	F	0.61	0/1031	0.82	0/1392
1	G	0.66	0/1040	0.82	0/1404
1	H	0.59	0/1095	0.85	0/1474
2	I	0.67	1/575 (0.2%)	0.89	0/887
2	K	0.81	2/557 (0.4%)	0.90	1/860 (0.1%)
2	M	0.88	4/575 (0.7%)	0.91	0/887
2	O	0.84	3/575 (0.5%)	0.92	1/887 (0.1%)
3	J	0.71	2/567 (0.4%)	0.94	0/872
3	L	0.84	3/567 (0.5%)	0.95	0/872
3	N	0.72	2/567 (0.4%)	0.88	0/872
3	P	0.94	4/543 (0.7%)	0.91	1/835 (0.1%)
All	All	0.70	21/13167 (0.2%)	0.86	3/18631 (0.0%)

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	K	4	DC	O3'-P	-7.75	1.51	1.61
2	K	5	DC	O3'-P	-7.67	1.51	1.61
3	P	5	DC	O3'-P	-7.20	1.52	1.61
2	O	4	DC	O3'-P	-6.99	1.52	1.61
2	M	6	DC	O3'-P	-6.83	1.52	1.61
3	N	16	DT	O3'-P	-6.66	1.53	1.61
3	L	5	DC	O3'-P	-6.32	1.53	1.61
3	L	4	DC	O3'-P	-6.15	1.53	1.61
3	N	11	DG	O3'-P	-6.00	1.53	1.61
2	M	14	DG	O3'-P	-5.97	1.53	1.61

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	M	5	DC	O3'-P	-5.86	1.54	1.61
2	O	13	DG	O3'-P	-5.85	1.54	1.61
3	J	16	DT	O3'-P	-5.70	1.54	1.61
3	P	16	DT	O3'-P	-5.52	1.54	1.61
2	M	13	DG	O3'-P	-5.43	1.54	1.61
2	O	14	DG	O3'-P	-5.40	1.54	1.61
3	L	16	DT	O3'-P	-5.36	1.54	1.61
2	I	4	DC	O3'-P	-5.31	1.54	1.61
3	P	10	DA	O3'-P	-5.27	1.54	1.61
3	P	14	DA	O3'-P	-5.11	1.55	1.61
3	J	14	DA	O3'-P	-5.09	1.55	1.61

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	O	14	DG	C1'-O4'-C4'	-5.97	104.13	110.10
2	K	4	DC	C1'-O4'-C4'	-5.64	104.46	110.10
3	P	16	DT	O5'-P-OP2	-5.18	101.03	105.70

There are no chirality outliers.

There are no planarity outliers.

## 5.2 Too-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 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	1150	0	1131	7	0
1	B	1007	0	1006	4	0
1	C	1021	0	1022	8	0
1	D	1150	0	1131	7	0
1	E	1089	0	1081	7	0
1	F	1021	0	1022	6	0
1	G	1030	0	1028	3	0
1	H	1085	0	1079	6	0
2	I	513	0	284	6	0
2	K	497	0	271	4	0
2	M	513	0	284	6	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	O	513	0	284	2	0
3	J	506	0	282	7	0
3	L	506	0	282	8	0
3	N	506	0	282	6	0
3	P	485	0	271	6	0
4	A	3	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
4	D	3	0	0	0	0
4	E	3	0	0	0	0
4	F	1	0	0	0	0
4	G	2	0	0	0	0
4	H	2	0	0	0	0
All	All	12608	0	10740	67	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 (67) 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:G:66:LEU:HD13	1:G:83:LEU:HD23	1.35	1.09
1:A:37:ARG:HG3	2:I:6:DC:OP1	1.57	1.05
1:C:66:LEU:HD13	1:C:83:LEU:HD13	1.54	0.89
1:D:71:ASP:HB3	3:N:3:DA:H5''	1.63	0.79
1:G:66:LEU:HD13	1:G:83:LEU:CD2	2.15	0.75
1:C:18:ARG:NH1	2:K:6:DC:OP2	2.21	0.74
1:F:18:ARG:NH1	3:N:6:DC:OP2	2.24	0.70
1:G:18:ARG:NH1	2:O:6:DC:OP2	2.25	0.69
1:H:71:ASP:HB3	3:J:3:DA:H5''	1.77	0.66
3:P:25:DA:H2'	3:P:25:DA:N3	2.10	0.65
1:B:18:ARG:NH1	3:J:6:DC:OP2	2.29	0.64
1:E:71:ASP:HB3	2:K:3:DA:OP1	1.99	0.63
1:A:15:GLU:OE2	3:J:17:DA:N7	2.32	0.62
1:C:66:LEU:HD13	1:C:83:LEU:CD1	2.28	0.61
1:F:107:GLU:O	1:F:110:ARG:HG2	2.01	0.60
1:F:107:GLU:O	1:F:110:ARG:CG	2.49	0.60
1:D:71:ASP:HB3	3:N:3:DA:C5'	2.32	0.60
1:D:65:SER:HB2	1:D:83:LEU:HD11	1.85	0.59
1:H:65:SER:HB2	1:H:83:LEU:HD11	1.84	0.59
2:M:12:DT:H2'	2:M:13:DG:C8	2.37	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:94:ARG:NH1	1:A:141:VAL:O	2.35	0.58
3:L:2:DC:H2'	3:L:3:DA:OP2	2.06	0.56
1:C:66:LEU:HD13	1:C:83:LEU:HD22	1.87	0.56
1:E:54:ARG:NH1	1:E:58:MET:O	2.39	0.55
1:C:66:LEU:CD1	1:C:83:LEU:HD13	2.32	0.55
3:L:4:DC:H2'	3:L:5:DC:C6	2.42	0.55
2:I:4:DC:H2'	2:I:5:DC:C6	2.43	0.54
2:I:10:DA:C2	3:J:17:DA:C2	2.96	0.54
1:H:71:ASP:HB3	3:J:3:DA:C5'	2.38	0.54
3:P:2:DC:H2'	3:P:3:DA:OP2	2.08	0.54
1:E:124:GLN:HA	1:E:124:GLN:OE1	2.08	0.53
1:H:71:ASP:OD2	3:J:3:DA:OP1	2.26	0.53
1:H:18:ARG:NH1	3:P:6:DC:OP2	2.42	0.52
1:A:29:PRO:HG2	1:A:37:ARG:HB3	1.90	0.52
1:E:18:ARG:NH1	2:M:6:DC:OP2	2.43	0.50
1:B:110:ARG:HB3	1:D:88:ILE:HG21	1.94	0.49
1:F:107:GLU:O	1:F:110:ARG:HG3	2.14	0.48
2:M:10:DA:C2	3:N:17:DA:C2	3.02	0.48
1:A:18:ARG:NH1	2:I:6:DC:OP2	2.43	0.47
3:L:4:DC:H2'	3:L:5:DC:H6	1.79	0.47
3:P:5:DC:H2'	3:P:6:DC:H6	1.79	0.47
2:I:4:DC:H2'	2:I:5:DC:H6	1.80	0.46
3:J:2:DC:H2'	3:J:3:DA:OP2	2.16	0.46
1:E:27:PRO:HG3	2:K:2:DG:OP2	2.16	0.45
3:N:2:DC:H2'	3:N:3:DA:OP2	2.16	0.45
2:O:17:DA:H1'	2:O:18:DC:H5'	1.97	0.45
3:L:5:DC:H2'	3:L:6:DC:O5'	2.16	0.45
1:A:37:ARG:NH1	2:I:7:DT:OP2	2.48	0.45
1:D:18:ARG:NH1	3:L:6:DC:OP2	2.44	0.44
1:C:105:LEU:HD22	1:D:84:ILE:HG23	2.00	0.43
2:M:5:DC:H2'	2:M:6:DC:C6	2.53	0.43
3:P:5:DC:H2'	3:P:6:DC:C6	2.53	0.43
1:E:71:ASP:CB	2:K:3:DA:OP1	2.65	0.43
1:E:49:PHE:HA	1:F:125:LEU:O	2.18	0.42
3:L:10:DA:O5'	3:L:10:DA:H2'	2.19	0.42
3:L:5:DC:H2'	3:L:6:DC:H6	1.84	0.42
1:B:103:GLU:HG2	1:C:109:ARG:HH22	1.84	0.42
2:M:17:DA:H2'	2:M:18:DC:O5'	2.20	0.42
1:F:110:ARG:HG3	1:F:111:ARG:N	2.35	0.41
1:B:6:LEU:HD22	1:B:39:TYR:CG	2.55	0.41
1:A:18:ARG:HD2	1:A:37:ARG:HH22	1.86	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:L:5:DC:H2'	3:L:6:DC:C6	2.56	0.41
1:C:66:LEU:HD13	1:C:83:LEU:CD2	2.51	0.41
3:P:15:DC:H2'	3:P:16:DT:C6	2.55	0.41
1:H:6:LEU:HD22	1:H:39:TYR:CG	2.57	0.40
1:D:43:HIS:CE1	3:N:1:DA:H5''	2.56	0.40
2:M:5:DC:H2'	2:M:6:DC:H6	1.86	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	143/147 (97%)	138 (96%)	5 (4%)	0	100	100
1	B	122/147 (83%)	119 (98%)	3 (2%)	0	100	100
1	C	124/147 (84%)	121 (98%)	3 (2%)	0	100	100
1	D	143/147 (97%)	138 (96%)	5 (4%)	0	100	100
1	E	131/147 (89%)	128 (98%)	3 (2%)	0	100	100
1	F	124/147 (84%)	120 (97%)	4 (3%)	0	100	100
1	G	125/147 (85%)	122 (98%)	3 (2%)	0	100	100
1	H	131/147 (89%)	128 (98%)	3 (2%)	0	100	100
All	All	1043/1176 (89%)	1014 (97%)	29 (3%)	0	100	100

There are no Ramachandran outliers to report.

### 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	126/127 (99%)	121 (96%)	5 (4%)	31	65
1	B	111/127 (87%)	109 (98%)	2 (2%)	59	85
1	C	113/127 (89%)	113 (100%)	0	100	100
1	D	126/127 (99%)	122 (97%)	4 (3%)	39	73
1	E	118/127 (93%)	114 (97%)	4 (3%)	37	71
1	F	113/127 (89%)	111 (98%)	2 (2%)	59	85
1	G	114/127 (90%)	111 (97%)	3 (3%)	46	77
1	H	118/127 (93%)	115 (98%)	3 (2%)	47	78
All	All	939/1016 (92%)	916 (98%)	23 (2%)	49	79

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	72	SER
1	A	102	GLN
1	A	132	SER
1	A	144	SER
1	A	145	HIS
1	B	72	SER
1	B	83	LEU
1	D	102	GLN
1	D	132	SER
1	D	144	SER
1	D	145	HIS
1	E	124	GLN
1	E	132	SER
1	E	144	SER
1	E	145	HIS
1	F	72	SER
1	F	102	GLN
1	G	72	SER
1	G	132	SER
1	G	135	GLU
1	H	102	GLN
1	H	144	SER
1	H	145	HIS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no

such sidechains identified.

### 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 monosaccharides in this entry.

## 5.6 Ligand geometry ⓘ

Of 16 ligands modelled in this entry, 16 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 ⓘ

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, 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	145/147 (98%)	0.15	3 (2%) 63 61	39, 60, 97, 118	0
1	B	126/147 (85%)	-0.06	0 100 100	31, 48, 78, 95	0
1	C	128/147 (87%)	0.03	2 (1%) 72 71	33, 50, 87, 104	0
1	D	145/147 (98%)	0.20	2 (1%) 75 75	37, 62, 105, 122	0
1	E	137/147 (93%)	0.17	5 (3%) 42 37	41, 64, 91, 115	0
1	F	128/147 (87%)	0.10	1 (0%) 86 86	34, 53, 83, 103	0
1	G	129/147 (87%)	0.09	2 (1%) 72 71	34, 50, 82, 95	0
1	H	137/147 (93%)	0.13	1 (0%) 87 87	30, 56, 87, 102	0
2	I	25/25 (100%)	-0.60	0 100 100	43, 65, 97, 125	0
2	K	25/25 (100%)	-0.68	0 100 100	37, 66, 95, 103	0
2	M	25/25 (100%)	-0.52	0 100 100	45, 71, 108, 135	0
2	O	25/25 (100%)	-0.70	0 100 100	34, 63, 90, 92	0
3	J	25/25 (100%)	-0.55	0 100 100	34, 59, 101, 122	0
3	L	25/25 (100%)	-0.62	0 100 100	43, 65, 104, 115	0
3	N	25/25 (100%)	-0.56	0 100 100	39, 63, 112, 131	0
3	P	24/25 (96%)	-0.61	0 100 100	44, 59, 86, 104	0
All	All	1274/1376 (92%)	-0.01	16 (1%) 77 77	30, 58, 95, 135	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	111	ARG	3.3
1	C	125	LEU	3.2
1	E	86	GLU	2.9
1	A	135	GLU	2.9
1	A	69	LEU	2.6

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Mol	Chain	Res	Type	RSRZ
1	E	11	ASP	2.6
1	E	84	ILE	2.5
1	D	133	VAL	2.4
1	F	74	ASP	2.2
1	A	138	HIS	2.2
1	D	89	GLU	2.2
1	E	138	HIS	2.2
1	H	76	SER	2.2
1	G	57	ASP	2.2
1	E	91	VAL	2.2
1	G	33	ASP	2.1

## 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 monosaccharides 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, 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	B-factors( $\text{\AA}^2$ )	Q<0.9
4	ZN	A	1001	1/1	0.94	0.11	91,91,91,91	0
4	ZN	D	202	1/1	0.96	0.10	101,101,101,101	0
4	ZN	G	202	1/1	0.98	0.13	91,91,91,91	0
4	ZN	D	203	1/1	0.99	0.16	51,51,51,51	0
4	ZN	B	201	1/1	0.99	0.17	38,38,38,38	0
4	ZN	G	201	1/1	0.99	0.16	40,40,40,40	0
4	ZN	D	201	1/1	0.99	0.13	59,59,59,59	0
4	ZN	H	201	1/1	0.99	0.18	44,44,44,44	0
4	ZN	E	202	1/1	0.99	0.11	118,118,118,118	0
4	ZN	E	201	1/1	1.00	0.15	50,50,50,50	0
4	ZN	H	202	1/1	1.00	0.14	64,64,64,64	0
4	ZN	F	201	1/1	1.00	0.14	45,45,45,45	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
4	ZN	E	203	1/1	1.00	0.14	63,63,63,63	0
4	ZN	A	1003	1/1	1.00	0.14	59,59,59,59	0
4	ZN	C	201	1/1	1.00	0.15	41,41,41,41	0
4	ZN	A	1002	1/1	1.00	0.17	50,50,50,50	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.