Full wwPDB X-ray Structure Validation Report (i)

Jan 29, 2024 – 04:29 PM EST

PDB ID	:	1F2I
Title	:	COCRYSTAL STRUCTURE OF SELECTED ZINC FINGER DIMER BOUND TO DNA
Authors	:	Wang, B.S.; Grant, R.A.; Pabo, C.O.
Deposited on	:	2000-05-25
Resolution	:	2.35 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 (i) 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 (i)) were used in the production of this report:

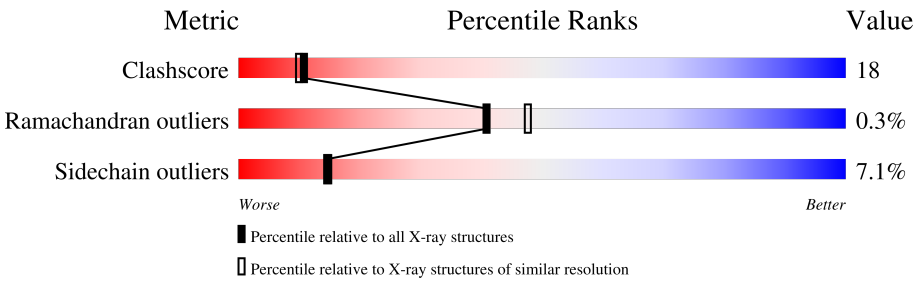
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

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 2.35 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))
Clashscore	141614	1232 (2.36-2.36)
Ramachandran outliers	138981	1211 (2.36-2.36)
Sidechain outliers	138945	1212 (2.36-2.36)

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 <=5%

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	14	64% 36%
1	B	14	71% 29%
1	C	14	71% 29%
1	D	14	64% 29% 7%
1	E	14	36% 57% 7%
1	F	14	64% 36%
2	G	73	60% 26% • 10%
2	H	73	59% 32% 10%

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Mol	Chain	Length	Quality of chain
2	I	73	55% 30% 5% 10%
2	J	73	60% 27% • 10%
2	K	73	49% 32% 5% 14%
2	L	73	38% 48% • 10%

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 5329 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 DNA chain called 5'-D(*AP*TP*GP*GP*GP*CP*GP*CP*GP*CP*CP*C P*AP*T)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
1	A	14	284	135	54	82	13	0	0	0
1	B	14	284	135	54	82	13	0	0	0
1	C	14	284	135	54	82	13	0	0	0
1	D	14	284	135	54	82	13	0	0	0
1	E	14	284	135	54	82	13	0	0	0
1	F	14	284	135	54	82	13	0	0	0

- Molecule 2 is a protein called FUSION OF N-TERMINAL 17-MER PEPTIDE EXTENSION TO ZIF12.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	G	66	553	340	113	94	6	0	0	0
2	H	66	553	340	113	94	6	0	0	0
2	I	66	553	340	113	94	6	0	0	0
2	J	66	553	340	113	94	6	0	0	0
2	K	63	529	324	109	90	6	0	0	0
2	L	66	553	340	113	94	6	0	0	0

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

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	G	2	Total Zn 2 2	0	0
3	H	2	Total Zn 2 2	0	0
3	I	2	Total Zn 2 2	0	0
3	J	2	Total Zn 2 2	0	0
3	K	2	Total Zn 2 2	0	0
3	L	2	Total Zn 2 2	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	22	Total O 22 22	0	0
4	B	15	Total O 15 15	0	0
4	C	25	Total O 25 25	0	0
4	D	35	Total O 35 35	0	0
4	E	10	Total O 10 10	0	0
4	F	8	Total O 8 8	0	0
4	G	35	Total O 35 35	0	0
4	H	39	Total O 39 39	0	0
4	I	42	Total O 42 42	0	0
4	J	42	Total O 42 42	0	0
4	K	26	Total O 26 26	0	0
4	L	20	Total O 20 20	0	0

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. 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. 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.

Note EDS was not executed.

• Molecule 1: 5'-D(*AP*TP*GP*GP*GP*CP*GP*CP*GP*CP*CP*CP*AP*T)-3'

Chain A: 64% 36%

• Molecule 1: 5'-D(*AP*TP*GP*GP*GP*CP*GP*CP*GP*CP*CP*CP*AP*T)-3'

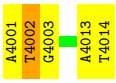
Chain B: 71% 29%

• Molecule 1: 5'-D(*AP*TP*GP*GP*GP*CP*GP*CP*GP*CP*CP*CP*AP*T)-3'

Chain C: 71% 29%

• Molecule 1: 5'-D(*AP*TP*GP*GP*GP*CP*GP*CP*GP*CP*CP*CP*AP*T)-3'

Chain D: 64% 29% 7%



• Molecule 1: 5'-D(*AP*TP*GP*GP*GP*CP*GP*CP*GP*CP*CP*CP*AP*T)-3'

Chain E: 36% 57% 7%

• Molecule 1: 5'-D(*AP*TP*GP*GP*GP*CP*GP*CP*GP*CP*CP*CP*AP*T)-3'

Chain F:

• Molecule 2: FUSION OF N-TERMINAL 17-MER PEPTIDE EXTENSION TO ZIF12

Chain G:

• Molecule 2: FUSION OF N-TERMINAL 17-MER PEPTIDE EXTENSION TO ZIF12

Chain H:

• Molecule 2: FUSION OF N-TERMINAL 17-MER PEPTIDE EXTENSION TO ZIF12

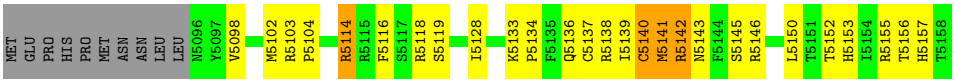
Chain I:

• Molecule 2: FUSION OF N-TERMINAL 17-MER PEPTIDE EXTENSION TO ZIF12

Chain J:

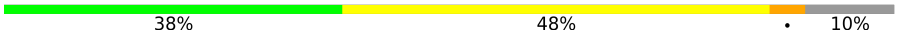
• Molecule 2: FUSION OF N-TERMINAL 17-MER PEPTIDE EXTENSION TO ZIF12

Chain K:



• Molecule 2: FUSION OF N-TERMINAL 17-MER PEPTIDE EXTENSION TO ZIF12

Chain L:



4 Data and refinement statistics (i)

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 31	Depositor
Cell constants a, b, c, α, β, γ	86.30Å 86.30Å 133.00Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 2.35	Depositor
% Data completeness (in resolution range)	82.6 (20.00-2.35)	Depositor
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR 3.851	Depositor
R, R_{free}	0.210 , 0.256	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	5329	wwPDB-VP
Average B, all atoms (Å^2)	41.0	wwPDB-VP

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	0.47	0/318	0.82	0/489
1	B	0.51	0/318	0.84	0/489
1	C	0.52	0/318	0.82	0/489
1	D	0.53	0/318	0.82	0/489
1	E	0.47	0/318	0.79	0/489
1	F	0.43	0/318	0.74	0/489
2	G	0.34	0/566	0.58	0/762
2	H	0.35	0/566	0.58	0/762
2	I	0.40	0/566	0.62	0/762
2	J	0.38	0/566	0.59	0/762
2	K	0.41	0/542	0.61	0/729
2	L	0.40	0/566	0.64	0/762
All	All	0.42	0/5280	0.69	0/7473

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand.A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	C	0	1
1	D	0	1
1	E	0	2
All	All	0	4

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	3009	DG	Sidechain
1	D	4002	DT	Sidechain
1	E	5003	DG	Sidechain
1	E	5009	DG	Sidechain

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	284	0	158	5	0
1	B	284	0	158	7	0
1	C	284	0	158	3	0
1	D	284	0	158	7	0
1	E	284	0	158	11	0
1	F	284	0	158	9	0
2	G	553	0	545	19	0
2	H	553	0	545	19	0
2	I	553	0	545	26	0
2	J	553	0	545	18	0
2	K	529	0	518	24	0
2	L	553	0	546	40	0
3	G	2	0	0	0	0
3	H	2	0	0	0	0
3	I	2	0	0	0	0
3	J	2	0	0	0	0
3	K	2	0	0	0	0
3	L	2	0	0	0	0
4	A	22	0	0	0	0
4	B	15	0	0	0	0
4	C	25	0	0	0	0
4	D	35	0	0	0	0
4	E	10	0	0	0	0
4	F	8	0	0	0	0
4	G	35	0	0	0	0
4	H	39	0	0	3	0
4	I	42	0	0	5	0
4	J	42	0	0	1	0
4	K	26	0	0	3	0
4	L	20	0	0	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	5329	0	4192	168	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 18.

All (168) 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)
2:I:3130:THR:HB	2:J:4095:LEU:HD23	1.26	1.11
1:F:6013:DA:H2"	1:F:6014:DT:H5"	1.31	1.07
1:A:1001:DA:H2"	1:A:1002:DT:H5'	1.41	1.02
2:K:5136:GLN:HG2	4:K:134:HOH:O	1.60	1.01
1:D:4001:DA:H2"	1:D:4002:DT:H5'	1.43	0.98
2:I:3130:THR:HB	2:J:4095:LEU:CD2	1.97	0.93
2:K:5141:MET:HA	4:K:134:HOH:O	1.73	0.88
1:B:2013:DA:H2"	1:B:2014:DT:H5"	1.59	0.85
2:K:5152:THR:O	2:K:5155:ARG:HG2	1.78	0.84
1:F:6004:DG:H5"	2:L:6114:ARG:HH21	1.41	0.82
2:I:3142:ARG:HH11	2:I:3142:ARG:HG2	1.46	0.81
2:G:1124:ARG:HD2	2:G:1146:ARG:HD2	1.65	0.79
2:L:6109:VAL:HG12	2:L:6112:CYS:H	1.47	0.78
2:H:2138:ARG:HG2	4:H:97:HOH:O	1.83	0.76
1:F:6013:DA:C2'	1:F:6014:DT:H5"	2.12	0.76
2:K:5136:GLN:HE21	2:K:5141:MET:HG3	1.55	0.71
1:F:6002:DT:H2"	1:F:6003:DG:C8	2.26	0.70
1:B:2004:DG:H2"	1:B:2005:DG:H5'	1.73	0.70
2:L:6110:GLU:HG3	2:L:6111:SER:H	1.56	0.70
1:D:4001:DA:H2"	1:D:4002:DT:C5'	2.19	0.69
1:F:6002:DT:H2"	1:F:6003:DG:H8	1.59	0.68
1:A:1001:DA:C2'	1:A:1002:DT:H5'	2.21	0.68
2:L:6146:ARG:HD2	4:L:88:HOH:O	1.92	0.68
1:B:2013:DA:H2"	1:B:2014:DT:C5'	2.23	0.68
1:B:2013:DA:C2'	1:B:2014:DT:H5"	2.24	0.67
2:K:5146:ARG:HD2	4:K:184:HOH:O	1.96	0.66
2:H:2099:VAL:HG23	4:H:91:HOH:O	1.97	0.64
2:G:1105:TYR:CE1	2:G:1119:SER:HB2	2.32	0.64
2:I:3130:THR:CB	2:J:4095:LEU:HD23	2.17	0.64
2:G:1095:LEU:CD2	2:H:2132:GLN:HE21	2.13	0.62
2:H:2122:LEU:O	2:H:2126:ILE:HG13	2.01	0.61
1:E:5001:DA:H2"	1:E:5002:DT:H5"	1.83	0.61
2:G:1143:ASN:HD22	2:G:1143:ASN:N	1.99	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:K:5137:CYS:HB2	2:K:5150:LEU:HD11	1.83	0.61
2:K:5136:GLN:NE2	2:K:5141:MET:HG3	2.16	0.60
2:L:6109:VAL:HG21	2:L:6129:HIS:CD2	2.36	0.60
2:G:1134:PRO:HG2	2:G:1135:PHE:CE1	2.36	0.60
2:I:3118:ARG:HB3	2:I:3120:ASP:OD1	2.01	0.60
1:F:6003:DG:H3'	2:L:6128:ILE:HD11	1.83	0.60
2:H:2140:CYS:O	2:H:2141:MET:HB2	2.01	0.59
2:L:6106:ALA:HB2	2:L:6115:ARG:HE	1.65	0.59
2:H:2142:ARG:HG2	2:H:2144:PHE:CZ	2.37	0.59
2:L:6150:LEU:O	2:L:6154:ILE:HG13	2.03	0.59
1:D:4002:DT:H1'	1:D:4003:DG:C8	2.38	0.58
2:K:5152:THR:O	2:K:5155:ARG:CG	2.51	0.58
1:F:6013:DA:H2"	1:F:6014:DT:C5'	2.20	0.57
2:J:4125:HIS:O	2:J:4128:ILE:HB	2.05	0.57
2:I:3106:ALA:HB2	2:I:3115:ARG:NH1	2.21	0.56
1:E:5003:DG:H3'	2:K:5128:ILE:HD11	1.88	0.56
1:E:5001:DA:H2"	1:E:5002:DT:C5'	2.36	0.56
2:L:6118:ARG:HB3	2:L:6120:ASP:OD1	2.06	0.55
1:E:5001:DA:C2'	1:E:5002:DT:H5"	2.36	0.55
1:C:3005:DG:OP1	2:I:3103:ARG:NH2	2.40	0.55
2:I:3107:CYS:HA	2:I:3122:LEU:HD11	1.89	0.55
2:I:3101:LYS:HD3	4:I:15:HOH:O	2.07	0.54
2:L:6093:ASN:OD1	2:L:6095:LEU:HB2	2.08	0.54
2:J:4103:ARG:HB3	2:J:4115:ARG:HB3	1.90	0.54
2:K:5139:ILE:HD12	2:K:5157:HIS:CD2	2.43	0.53
2:I:3142:ARG:HH11	2:I:3142:ARG:CG	2.18	0.53
2:J:4130:THR:HB	2:J:4132:GLN:HG3	1.90	0.53
2:H:2116:PHE:HB3	2:H:2121:GLU:OE1	2.08	0.53
2:I:3102:MET:HB2	4:I:288:HOH:O	2.08	0.53
2:L:6158:THR:HG22	2:L:6158:THR:O	2.08	0.53
1:E:5005:DG:OP1	2:K:5103:ARG:NH2	2.42	0.52
2:L:6098:VAL:HA	4:L:177:HOH:O	2.09	0.52
1:E:5011:DC:H5'	1:E:5011:DC:H6	1.73	0.52
2:G:1095:LEU:HD22	2:H:2132:GLN:HE21	1.73	0.52
2:G:1102:MET:HE1	2:G:1104:PRO:HG3	1.91	0.52
2:K:5133:LYS:N	2:K:5134:PRO:HD3	2.25	0.52
2:I:3103:ARG:HB2	2:I:3115:ARG:HB3	1.91	0.52
1:F:6003:DG:C5'	2:L:6128:ILE:HD11	2.40	0.52
2:I:3142:ARG:HG2	2:I:3142:ARG:NH1	2.22	0.52
2:J:4100:PRO:O	2:J:4102:MET:HG3	2.09	0.52
1:B:2014:DT:H6	1:B:2014:DT:H5'	1.75	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:L:6122:LEU:O	2:L:6126:ILE:HG13	2.10	0.52
2:K:5103:ARG:HA	2:K:5116:PHE:O	2.10	0.51
2:L:6140:CYS:O	2:L:6141:MET:HB2	2.11	0.51
2:L:6135:PHE:O	2:L:6144:PHE:N	2.44	0.51
2:G:1118:ARG:HB3	2:G:1120:ASP:OD1	2.11	0.51
2:L:6154:ILE:HG22	2:L:6154:ILE:O	2.11	0.50
2:I:3101:LYS:HA	4:I:15:HOH:O	2.12	0.50
2:L:6103:ARG:HB2	2:L:6115:ARG:HB3	1.93	0.50
2:L:6153:HIS:C	2:L:6155:ARG:H	2.15	0.49
1:D:4003:DG:H5”	2:J:4128:ILE:HD11	1.94	0.49
2:G:1135:PHE:HB3	2:G:1150:LEU:HD22	1.94	0.49
2:G:1118:ARG:HG2	2:H:2118:ARG:HG2	1.93	0.49
2:G:1137:CYS:SG	2:G:1139:ILE:HB	2.52	0.49
2:L:6103:ARG:CB	2:L:6115:ARG:HB3	2.43	0.49
2:L:6109:VAL:HG11	2:L:6112:CYS:HB2	1.95	0.49
2:K:5140:CYS:SG	2:K:5142:ARG:HB2	2.52	0.49
2:I:3114:ARG:HG3	2:I:3125:HIS:CE1	2.48	0.48
2:I:3093:ASN:C	2:I:3095:LEU:H	2.16	0.48
1:D:4003:DG:H3’	2:J:4128:ILE:HD11	1.95	0.48
2:H:2126:ILE:HG22	2:H:2126:ILE:O	2.14	0.48
2:J:4142:ARG:HH11	2:J:4153:HIS:HE1	1.61	0.48
2:I:3103:ARG:CB	2:I:3115:ARG:HB3	2.44	0.48
2:L:6122:LEU:HG	2:L:6126:ILE:HD11	1.95	0.47
2:H:2109:VAL:HG21	2:H:2129:HIS:CD2	2.49	0.47
2:L:6102:MET:O	2:L:6104:PRO:HD3	2.15	0.47
2:L:6101:LYS:O	2:L:6101:LYS:HG2	2.15	0.47
2:L:6109:VAL:CG1	2:L:6112:CYS:HB2	2.45	0.47
2:K:5133:LYS:HA	2:K:5145:SER:HA	1.95	0.47
1:A:1003:DG:H3’	2:G:1128:ILE:HD11	1.97	0.47
2:G:1118:ARG:HG3	4:H:22:HOH:O	2.14	0.46
2:L:6140:CYS:SG	2:L:6142:ARG:HB2	2.55	0.46
2:K:5114:ARG:HH11	2:K:5114:ARG:HG3	1.80	0.46
2:L:6112:CYS:SG	2:L:6114:ARG:HB2	2.56	0.46
2:G:1120:ASP:OD1	2:G:1120:ASP:N	2.46	0.46
2:H:2102:MET:C	2:H:2103:ARG:HG3	2.35	0.46
2:K:5114:ARG:HG2	2:K:5116:PHE:CZ	2.51	0.46
2:J:4118:ARG:HB3	2:J:4120:ASP:OD1	2.16	0.45
2:H:2118:ARG:HB3	2:H:2120:ASP:OD1	2.16	0.45
2:L:6149:HIS:HD2	4:L:174:HOH:O	1.99	0.45
2:L:6146:ARG:HB3	2:L:6148:ASP:OD1	2.15	0.45
2:L:6141:MET:HG2	4:L:50:HOH:O	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:L:6156:THR:O	2:L:6156:THR:HG22	2.17	0.45
1:C:3013:DA:H2"	1:C:3014:DT:H5'	2.00	0.44
2:G:1143:ASN:N	2:G:1143:ASN:ND2	2.64	0.44
2:H:2103:ARG:O	2:H:2115:ARG:NH1	2.48	0.44
1:B:2005:DG:OP1	2:H:2103:ARG:NH2	2.46	0.44
1:D:4013:DA:H2"	1:D:4014:DT:C5'	2.48	0.44
2:G:1137:CYS:HA	2:G:1154:ILE:HD11	1.98	0.44
2:I:3135:PHE:HB3	2:I:3150:LEU:HD22	2.00	0.44
1:D:4013:DA:H2"	1:D:4014:DT:H5'	2.00	0.44
2:K:5098:VAL:O	2:K:5098:VAL:HG23	2.18	0.44
1:C:3013:DA:H2"	1:C:3014:DT:C5'	2.48	0.44
2:L:6157:HIS:O	2:L:6158:THR:C	2.56	0.43
2:I:3111:SER:HB3	4:I:201:HOH:O	2.18	0.43
2:K:5114:ARG:HG3	2:K:5114:ARG:NH1	2.32	0.43
2:I:3120:ASP:OD1	2:I:3120:ASP:N	2.51	0.43
2:J:4138:ARG:HE	2:J:4138:ARG:HB2	1.62	0.43
2:L:6135:PHE:HB3	2:L:6150:LEU:HD23	2.01	0.43
4:I:99:HOH:O	2:J:4099:VAL:HG23	2.17	0.43
2:H:2107:CYS:HA	2:H:2122:LEU:HD11	2.01	0.43
2:I:3098:VAL:O	2:I:3100:PRO:HD3	2.18	0.43
2:I:3142:ARG:CG	2:I:3142:ARG:NH1	2.78	0.42
2:L:6150:LEU:O	2:L:6150:LEU:HD12	2.18	0.42
1:E:5005:DG:H1'	1:E:5006:DC:H5'	2.01	0.42
1:E:5014:DT:H6	1:E:5014:DT:H5'	1.85	0.42
2:G:1093:ASN:HD22	2:G:1093:ASN:HA	1.65	0.42
2:L:6107:CYS:C	2:L:6109:VAL:H	2.22	0.42
2:K:5156:THR:O	2:K:5156:THR:HG22	2.19	0.42
2:H:2142:ARG:HG2	2:H:2144:PHE:CE1	2.54	0.42
2:G:1124:ARG:O	2:G:1127:ARG:HB2	2.20	0.42
2:J:4140:CYS:O	2:J:4141:MET:HB2	2.20	0.42
1:B:2004:DG:H2"	1:B:2005:DG:C5'	2.47	0.41
1:E:5001:DA:H1'	1:E:5002:DT:H5"	2.01	0.41
2:J:4099:VAL:HA	2:J:4100:PRO:HD2	1.91	0.41
1:A:1003:DG:N7	2:G:1149:HIS:NE2	2.57	0.41
1:E:5010:DC:H2"	1:E:5011:DC:H5'	2.03	0.41
2:L:6126:ILE:HG22	2:L:6126:ILE:O	2.20	0.41
1:A:1007:DG:H2"	1:A:1008:DC:C6	2.55	0.41
1:E:5010:DC:H2"	1:E:5011:DC:C5'	2.51	0.41
2:I:3093:ASN:C	2:I:3095:LEU:N	2.74	0.41
2:K:5118:ARG:HG2	2:L:6118:ARG:HG2	2.02	0.41
2:L:6140:CYS:O	2:L:6141:MET:CB	2.69	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:I:3124:ARG:HD2	2:I:3146:ARG:HG2	2.03	0.40
2:K:5136:GLN:HG3	2:K:5143:ASN:OD1	2.21	0.40
2:H:2105:TYR:CE1	2:H:2119:SER:HB3	2.57	0.40
2:I:3118:ARG:HG2	2:J:4118:ARG:HG2	2.04	0.40
2:J:4109:VAL:HB	2:J:4112:CYS:HB3	2.03	0.40
2:K:5102:MET:HE1	2:K:5104:PRO:HG3	2.02	0.40
2:L:6152:THR:O	2:L:6152:THR:HG22	2.21	0.40
2:H:2098:VAL:HG23	2:H:2098:VAL:O	2.22	0.40
2:I:3094:LEU:HA	2:I:3097:TYR:HB2	2.04	0.40
2:J:4138:ARG:HD3	4:J:292:HOH:O	2.20	0.40
2:K:5155:ARG:C	2:K:5157:HIS:N	2.75	0.40
1:F:6003:DG:H5"	2:L:6128:ILE:HD11	2.02	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	
2	G	64/73 (88%)	62 (97%)	2 (3%)	0	100	100
2	H	64/73 (88%)	64 (100%)	0	0	100	100
2	I	64/73 (88%)	60 (94%)	3 (5%)	1 (2%)	9	8
2	J	64/73 (88%)	61 (95%)	3 (5%)	0	100	100
2	K	61/73 (84%)	52 (85%)	9 (15%)	0	100	100
2	L	64/73 (88%)	57 (89%)	7 (11%)	0	100	100
All	All	381/438 (87%)	356 (93%)	24 (6%)	1 (0%)	41	47

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	I	3100	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	
2	G	64/71 (90%)	59 (92%)	5 (8%)	12	12
2	H	64/71 (90%)	63 (98%)	1 (2%)	62	75
2	I	64/71 (90%)	59 (92%)	5 (8%)	12	12
2	J	64/71 (90%)	59 (92%)	5 (8%)	12	12
2	K	61/71 (86%)	54 (88%)	7 (12%)	5	5
2	L	64/71 (90%)	60 (94%)	4 (6%)	18	19
All	All	381/426 (89%)	354 (93%)	27 (7%)	14	15

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	G	1093	ASN
2	G	1098	VAL
2	G	1120	ASP
2	G	1141	MET
2	G	1143	ASN
2	H	2147	SER
2	I	3093	ASN
2	I	3102	MET
2	I	3114	ARG
2	I	3138	ARG
2	I	3147	SER
2	J	4114	ARG
2	J	4117	SER
2	J	4119	SER
2	J	4128	ILE
2	J	4138	ARG
2	K	5114	ARG
2	K	5119	SER
2	K	5138	ARG
2	K	5140	CYS
2	K	5141	MET
2	K	5142	ARG

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Mol	Chain	Res	Type
2	K	5153	HIS
2	L	6114	ARG
2	L	6117	SER
2	L	6140	CYS
2	L	6141	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
2	G	1093	ASN
2	G	1132	GLN
2	G	1143	ASN
2	H	2132	GLN
2	I	3093	ASN
2	J	4136	GLN
2	K	5136	GLN
2	K	5153	HIS
2	K	5157	HIS

5.3.3 RNA (i)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains (i)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 12 ligands modelled in this entry, 12 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)

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains (i)

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates (i)

EDS was not executed - this section is therefore empty.

6.4 Ligands (i)

EDS was not executed - this section is therefore empty.

6.5 Other polymers (i)

EDS was not executed - this section is therefore empty.