



# Full wwPDB X-ray Structure Validation Report ⓘ

May 22, 2020 – 07:23 pm BST

PDB ID : 1AIS  
Title : TATA-BINDING PROTEIN/TRANSCRIPTION FACTOR (II)B/TATA-BOX COMPLEX FROM PYROCOCCLUS WOESEI  
Authors : Kosa, P.F.; Ghosh, G.; Dedecker, B.S.; Sigler, P.B.  
Deposited on : 1997-04-24  
Resolution : 2.10 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  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
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

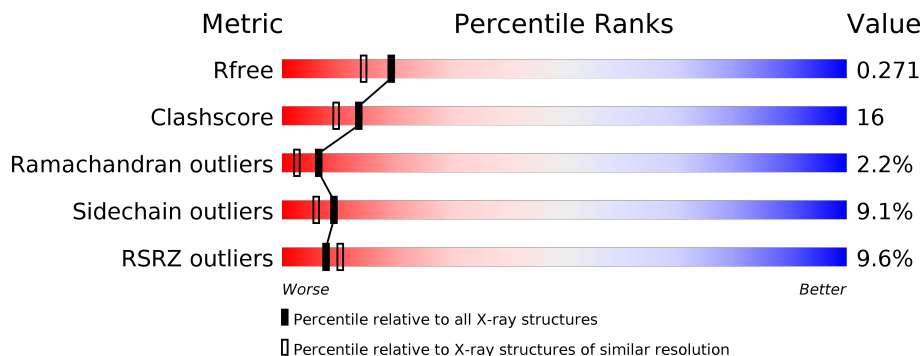
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.10 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	5197 (2.10-2.10)
Clashscore	141614	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (2.10-2.10)
RSRZ outliers	127900	5083 (2.10-2.10)

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	C	17	
2	E	17	
3	A	182	
4	B	200	

## 2 Entry composition i

There are 5 unique types of molecules in this entry. The entry contains 3929 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 DNA (5'-D(\*AP\*AP\*CP\*TP\*TP\*AP\*CP\*TP\*TP\*TP\*(5IU)P\*(5IU)P\*AP\*AP\*AP\*GP\*C)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace	
			Total	C	I	N	O				P
1	C	17	342	165	2	58	101	16	0	0	0

- Molecule 2 is a DNA chain called DNA (5'-D(\*GP\*CP\*TP\*TP\*TP\*AP\*AP\*AP\*AP\*AP\*GP\*TP\*AP\*AP\*GP\*TP\*T)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
2	E	17	349	169	65	99	16	0	0	0

- Molecule 3 is a protein called PROTEIN (TATA-BINDING PROTEIN).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	A	181	1414	910	235	262	7	0	0	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	182	LEU	-	INSERTION	UNP P62001

- Molecule 4 is a protein called PROTEIN (TRANSCRIPTION INITIATION FACTOR IIB).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
4	B	193	1536	972	285	276	3	0	0	0

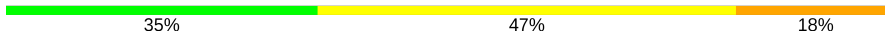
- Molecule 5 is water.

<b>Mol</b>	<b>Chain</b>	<b>Residues</b>	<b>Atoms</b>		<b>ZeroOcc</b>	<b>AltConf</b>
5	C	36	Total 36	O 36	1	0
5	E	45	Total 45	O 45	0	0
5	A	146	Total 146	O 146	0	0
5	B	61	Total 61	O 61	0	0

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

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.

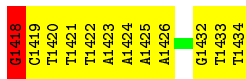
- Molecule 1: DNA (5'-D(\*AP\*AP\*CP\*TP\*TP\*AP\*CP\*TP\*TP\*TP\*(5IU)P\*(5IU)P\*AP\*AP\*AP\*GP\*C)-3')

Chain C: 




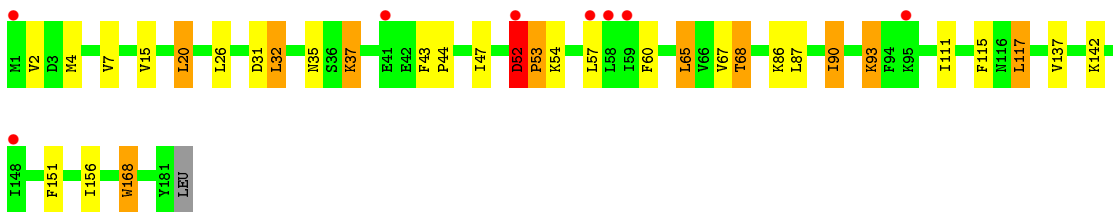
- Molecule 2: DNA (5'-D(\*GP\*CP\*TP\*TP\*TP\*AP\*AP\*AP\*AP\*AP\*GP\*TP\*AP\*AP\*GP\*TP\*T)-3')

Chain E: 



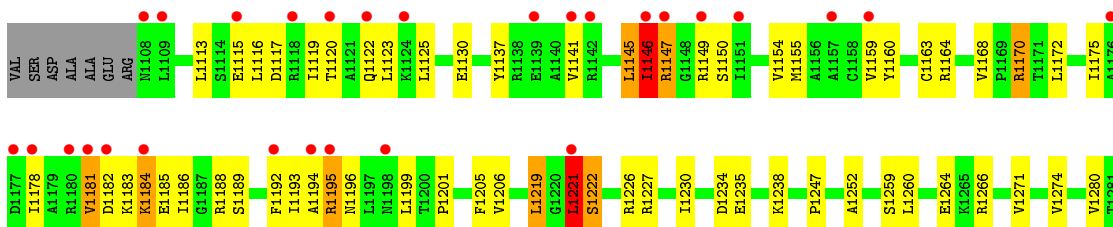
- Molecule 3: PROTEIN (TATA-BINDING PROTEIN)

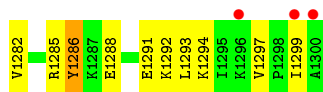
Chain A: 



- Molecule 4: PROTEIN (TRANSCRIPTION INITIATION FACTOR IIB)

Chain B: 





## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	125.70Å 91.20Å 74.20Å 90.00° 122.70° 90.00°	Depositor
Resolution (Å)	20.00 – 2.10 24.78 – 2.08	Depositor EDS
% Data completeness (in resolution range)	92.0 (20.00-2.10) 96.1 (24.78-2.08)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.09	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.10 (at 2.08Å)	Xtrriage
Refinement program	X-PLOR	Depositor
R, $R_{free}$	0.212 , 0.268 0.217 , 0.271	Depositor DCC
$R_{free}$ test set	2009 reflections (4.96%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	31.4	Xtrriage
Anisotropy	0.286	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.39 , 78.7	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	3929	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	36.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.83% 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 [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: 5IU

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	C	0.59	0/337	1.04	1/515 (0.2%)
2	E	0.60	0/392	1.03	1/604 (0.2%)
3	A	0.50	0/1437	0.75	2/1939 (0.1%)
4	B	0.43	0/1553	0.65	1/2086 (0.0%)
All	All	0.50	0/3719	0.79	5/5144 (0.1%)

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	2
2	E	0	3
All	All	0	5

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
3	A	52	ASP	C-N-CD	6.26	141.54	128.40
2	E	1418	DG	N9-C1'-C2'	-5.46	102.23	112.60
4	B	1221	LEU	N-CA-C	5.36	125.47	111.00
1	C	1401	DA	N9-C1'-C2'	-5.22	102.68	112.60
3	A	53	PRO	N-CA-C	-5.15	98.70	112.10

There are no chirality outliers.

All (5) planarity outliers are listed below:



Mol	Chain	Res	Type	Group
1	C	1407	DC	Sidechain
1	C	1414	DA	Sidechain
2	E	1418	DG	Sidechain
2	E	1424	DA	Sidechain
2	E	1425	DA	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	C	342	0	190	22	0
2	E	349	0	195	18	0
3	A	1414	0	1473	22	0
4	B	1536	0	1639	55	0
5	A	146	0	0	0	0
5	B	61	0	0	2	0
5	C	36	0	0	2	0
5	E	45	0	0	0	0
All	All	3929	0	3497	113	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 16.

All (113) 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:E:1418:DG:H2''	2:E:1419:DC:C5'	1.72	1.19
2:E:1418:DG:H2''	2:E:1419:DC:H5''	1.29	1.08
1:C:1403:DC:H2''	1:C:1404:DT:H5'	1.14	1.08
2:E:1418:DG:H5'	2:E:1418:DG:H8	1.29	0.97
2:E:1419:DC:H5'	2:E:1419:DC:H6	1.27	0.96
2:E:1421:DT:H2''	2:E:1422:DT:H5'	1.48	0.96
1:C:1403:DC:H2''	1:C:1404:DT:C5'	2.00	0.92
2:E:1418:DG:H2''	2:E:1419:DC:H5'	1.52	0.89
1:C:1402:DA:H2''	1:C:1403:DC:H5'	1.55	0.89
1:C:1402:DA:H2''	1:C:1403:DC:C5'	2.03	0.87
1:C:1403:DC:C2'	1:C:1404:DT:H5'	2.02	0.87

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:1418:DG:C8	2:E:1418:DG:H5'	2.13	0.83
2:E:1421:DT:H2''	2:E:1422:DT:C5'	2.10	0.81
3:A:52:ASP:HB3	3:A:53:PRO:HD3	1.61	0.81
2:E:1418:DG:C2'	2:E:1419:DC:C5'	2.60	0.77
2:E:1418:DG:C2'	2:E:1419:DC:H5'	2.15	0.77
3:A:37:LYS:HG2	3:A:47:ILE:HB	1.69	0.74
4:B:1149:ARG:NH1	4:B:1192:PHE:HB3	2.03	0.74
4:B:1221:LEU:HB3	4:B:1226:ARG:HH21	1.51	0.74
3:A:15:VAL:HG22	3:A:68:THR:HG22	1.71	0.72
3:A:4:MET:O	3:A:7:VAL:HG12	1.88	0.72
4:B:1146:ILE:HB	4:B:1149:ARG:HE	1.55	0.71
1:C:1407:DC:H2'	1:C:1408:DT:H72	1.73	0.68
3:A:117:LEU:HG	3:A:137:VAL:HG23	1.75	0.68
1:C:1412:5IU:I5	5:C:1677:HOH:O	2.82	0.68
4:B:1181:VAL:HG22	4:B:1182:ASP:H	1.59	0.67
3:A:15:VAL:HG22	3:A:68:THR:CG2	2.26	0.66
3:A:52:ASP:HB3	3:A:53:PRO:CD	2.26	0.64
4:B:1147:ARG:HB2	4:B:1147:ARG:HH11	1.63	0.64
4:B:1146:ILE:HD12	4:B:1149:ARG:NE	2.11	0.64
1:C:1402:DA:H2''	1:C:1403:DC:H5''	1.78	0.64
4:B:1145:LEU:HD21	4:B:1154:VAL:HG11	1.80	0.63
2:E:1419:DC:H5'	2:E:1419:DC:C6	2.20	0.63
3:A:20:LEU:HD22	3:A:65:LEU:HD21	1.81	0.63
3:A:32:LEU:HG	3:A:86:LYS:HB3	1.81	0.63
1:C:1407:DC:C2'	1:C:1408:DT:H72	2.28	0.62
4:B:1172:LEU:HD11	4:B:1186:ILE:HG21	1.81	0.62
4:B:1116:LEU:HD22	4:B:1130:GLU:HG3	1.83	0.60
3:A:111:ILE:HD11	3:A:115:PHE:CE1	2.36	0.59
4:B:1116:LEU:CD2	4:B:1130:GLU:HG3	2.32	0.59
4:B:1159:VAL:HG23	4:B:1175:ILE:HD13	1.84	0.59
1:C:1408:DT:H2''	1:C:1409:DT:H5'	1.85	0.59
4:B:1149:ARG:HH11	4:B:1189:SER:HA	1.68	0.59
4:B:1192:PHE:O	4:B:1196:ASN:HB2	2.02	0.59
2:E:1419:DC:H2'	2:E:1420:DT:H71	1.86	0.57
4:B:1146:ILE:HD12	4:B:1149:ARG:CD	2.33	0.57
1:C:1414:DA:H1'	3:A:60:PHE:CZ	2.40	0.57
4:B:1147:ARG:HB2	4:B:1147:ARG:NH1	2.20	0.57
4:B:1259:SER:HB2	4:B:1264:GLU:O	2.05	0.56
4:B:1234:ASP:O	4:B:1238:LYS:HG2	2.06	0.56
4:B:1221:LEU:CB	4:B:1226:ARG:HH21	2.19	0.56
4:B:1221:LEU:HB3	4:B:1226:ARG:NH2	2.19	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:1423:DA:N3	3:A:68:THR:HG21	2.21	0.56
4:B:1125:LEU:HD21	4:B:1170:ARG:HD2	1.90	0.54
4:B:1183:LYS:HD2	4:B:1183:LYS:N	2.22	0.54
4:B:1282:VAL:O	4:B:1286:TYR:HB3	2.07	0.54
4:B:1235:GLU:HG3	4:B:1293:LEU:HD22	1.89	0.54
2:E:1433:DT:H2'	2:E:1434:DT:H71	1.88	0.54
4:B:1146:ILE:HG22	4:B:1149:ARG:H	1.74	0.53
4:B:1117:ASP:HA	4:B:1120:THR:HG22	1.90	0.53
4:B:1260:LEU:HD13	4:B:1297:VAL:HG21	1.91	0.53
2:E:1418:DG:H2'	2:E:1419:DC:C6	2.44	0.52
4:B:1194:ALA:O	4:B:1199:LEU:HB2	2.09	0.52
1:C:1403:DC:C2'	1:C:1404:DT:C5'	2.77	0.51
4:B:1122:GLN:NE2	4:B:1178:ILE:HB	2.25	0.51
4:B:1183:LYS:H	4:B:1183:LYS:HD2	1.75	0.51
4:B:1195:ARG:HA	4:B:1195:ARG:NE	2.26	0.51
4:B:1155:MET:O	4:B:1159:VAL:HG13	2.10	0.51
3:A:111:ILE:HG23	3:A:156:ILE:HD13	1.94	0.50
4:B:1280:VAL:HG23	5:B:1562:HOH:O	2.11	0.50
1:C:1407:DC:C6	1:C:1408:DT:H72	2.47	0.50
3:A:57:LEU:HD23	3:A:67:VAL:HG22	1.94	0.50
1:C:1414:DA:H1'	3:A:60:PHE:CE2	2.47	0.50
4:B:1160:TYR:O	4:B:1164:ARG:HG3	2.12	0.49
1:C:1408:DT:C6	1:C:1409:DT:H72	2.47	0.49
4:B:1222:SER:O	4:B:1226:ARG:HG3	2.12	0.49
1:C:1402:DA:C2'	1:C:1403:DC:H5''	2.41	0.49
1:C:1411:5IU:I5	5:C:1636:HOH:O	2.90	0.48
4:B:1189:SER:O	4:B:1193:ILE:HG22	2.13	0.48
4:B:1163:CYS:HB3	4:B:1168:VAL:O	2.14	0.48
2:E:1426:DA:H1'	3:A:151:PHE:CE1	2.50	0.47
4:B:1146:ILE:HG22	4:B:1149:ARG:N	2.29	0.47
4:B:1291:GLU:HG3	4:B:1292:LYS:N	2.30	0.47
3:A:37:LYS:CG	3:A:47:ILE:HB	2.41	0.46
4:B:1172:LEU:HD11	4:B:1186:ILE:CG2	2.46	0.46
4:B:1137:TYR:O	4:B:1141:VAL:HG23	2.15	0.46
4:B:1116:LEU:HD23	4:B:1116:LEU:O	2.16	0.45
2:E:1432:DG:H2''	2:E:1433:DT:H5'	1.98	0.45
4:B:1288:GLU:HG3	4:B:1292:LYS:HE2	1.98	0.45
1:C:1402:DA:C2'	1:C:1403:DC:C5'	2.87	0.45
4:B:1227:ARG:HH12	4:B:1230:ILE:HG21	1.81	0.45
3:A:90:ILE:O	3:A:90:ILE:HG23	2.17	0.45
3:A:52:ASP:O	3:A:54:LYS:N	2.50	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:B:1188:ARG:O	4:B:1192:PHE:HB2	2.17	0.45
3:A:93:LYS:NZ	3:A:93:LYS:HB2	2.32	0.45
4:B:1123:LEU:HB2	4:B:1125:LEU:HG	2.00	0.44
4:B:1219:LEU:HD13	4:B:1274:VAL:HG11	1.99	0.44
4:B:1201:PRO:O	4:B:1205:PHE:HB2	2.18	0.44
1:C:1407:DC:C2'	1:C:1408:DT:C7	2.95	0.43
3:A:43:PHE:CD1	3:A:44:PRO:HD2	2.53	0.43
4:B:1291:GLU:HG3	4:B:1292:LYS:H	1.83	0.43
1:C:1403:DC:H6	1:C:1403:DC:H5'	1.82	0.43
4:B:1247:PRO:HD2	5:B:1547:HOH:O	2.18	0.42
4:B:1115:GLU:O	4:B:1119:ILE:HG13	2.19	0.42
4:B:1119:ILE:HG21	4:B:1159:VAL:HG11	2.02	0.42
4:B:1252:ALA:O	4:B:1271:VAL:HG11	2.19	0.42
4:B:1116:LEU:O	4:B:1120:THR:HG22	2.20	0.42
1:C:1410:DT:H2''	1:C:1411:5IU:H5''	2.00	0.42
1:C:1401:DA:H2'	1:C:1402:DA:C8	2.54	0.41
4:B:1145:LEU:CD2	4:B:1154:VAL:HG11	2.49	0.41
2:E:1433:DT:H5'	2:E:1433:DT:H6	1.85	0.41
3:A:7:VAL:HG13	3:A:168:TRP:CZ2	2.56	0.40
4:B:1294:LYS:HA	4:B:1294:LYS:HD3	1.96	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
3	A	179/182 (98%)	176 (98%)	2 (1%)	1 (1%)	25 21
4	B	191/200 (96%)	173 (91%)	11 (6%)	7 (4%)	3 1
All	All	370/382 (97%)	349 (94%)	13 (4%)	8 (2%)	6 2

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	A	52	ASP
4	B	1146	ILE
4	B	1181	VAL
4	B	1221	LEU
4	B	1299	ILE
4	B	1184	LYS
4	B	1222	SER
4	B	1150	SER

### 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
3	A	159/160 (99%)	144 (91%)	15 (9%)	8 5
4	B	161/166 (97%)	147 (91%)	14 (9%)	10 7
All	All	320/326 (98%)	291 (91%)	29 (9%)	9 6

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	A	2	VAL
3	A	20	LEU
3	A	26	LEU
3	A	31	ASP
3	A	32	LEU
3	A	35	ASN
3	A	37	LYS
3	A	65	LEU
3	A	68	THR
3	A	87	LEU
3	A	90	ILE
3	A	93	LYS
3	A	117	LEU
3	A	142	LYS
3	A	168	TRP
4	B	1113	LEU

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Mol	Chain	Res	Type
4	B	1145	LEU
4	B	1146	ILE
4	B	1147	ARG
4	B	1170	ARG
4	B	1184	LYS
4	B	1185	GLU
4	B	1195	ARG
4	B	1206	VAL
4	B	1219	LEU
4	B	1221	LEU
4	B	1266	ARG
4	B	1285	ARG
4	B	1286	TYR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
3	A	85	GLN
3	A	103	GLN
4	B	1122	GLN
4	B	1198	ASN
4	B	1213	ASN
4	B	1268	GLN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

2 non-standard protein/DNA/RNA residues are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
1	5IU	C	1411	1,2	14,21,22	1.49	2 (14%)	16,30,33	4.21	4 (25%)
1	5IU	C	1412	1,2	14,21,22	1.49	3 (21%)	16,30,33	4.21	3 (18%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	5IU	C	1411	1,2	-	1/4/21/22	0/2/2/2
1	5IU	C	1412	1,2	-	2/4/21/22	0/2/2/2

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	1411	5IU	C6-C5	-4.17	1.33	1.38
1	C	1412	5IU	C6-C5	-3.33	1.34	1.38
1	C	1411	5IU	C4-N3	3.20	1.38	1.33
1	C	1412	5IU	C4-N3	3.02	1.38	1.33
1	C	1412	5IU	C5-I5	2.84	2.16	2.10

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	1412	5IU	C4-N3-C2	14.32	127.23	115.14
1	C	1411	5IU	C4-N3-C2	14.16	127.10	115.14
1	C	1411	5IU	C5-C6-N1	6.05	124.10	120.70
1	C	1411	5IU	C5-C4-N3	-5.97	115.28	123.27
1	C	1412	5IU	C5-C4-N3	-5.97	115.28	123.27
1	C	1412	5IU	C5-C6-N1	5.75	123.94	120.70
1	C	1411	5IU	C2'-C1'-N1	-2.04	109.57	114.27

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	C	1411	5IU	O4'-C1'-N1-C6
1	C	1412	5IU	O4'-C1'-N1-C6
1	C	1412	5IU	C3'-C4'-C5'-O5'

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	C	1411	5IU	2	0
1	C	1412	5IU	1	0

## 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

There are no ligands in this entry.

## 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, 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	C	15/17 (88%)	-0.45	0 <span style="border: 1px solid blue; padding: 2px;">100</span> <span style="border: 1px solid blue; padding: 2px;">100</span>	16, 21, 42, 47	0
2	E	17/17 (100%)	-0.65	0 <span style="border: 1px solid blue; padding: 2px;">100</span> <span style="border: 1px solid blue; padding: 2px;">100</span>	13, 21, 42, 44	0
3	A	181/182 (99%)	-0.05	8 (4%) <span style="border: 1px solid red; padding: 2px;">34</span> <span style="border: 1px solid red; padding: 2px;">40</span>	10, 22, 47, 70	0
4	B	193/200 (96%)	0.73	31 (16%) <span style="border: 1px solid red; padding: 2px;">1</span> <span style="border: 1px solid red; padding: 2px;">2</span>	14, 44, 95, 100	0
All	All	406/416 (97%)	0.28	39 (9%) <span style="border: 1px solid red; padding: 2px;">8</span> <span style="border: 1px solid red; padding: 2px;">10</span>	10, 29, 91, 100	0

All (39) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
4	B	1147	ARG	10.6
4	B	1300	ALA	5.6
4	B	1146	ILE	5.4
4	B	1180	ARG	5.2
4	B	1182	ASP	4.8
4	B	1181	VAL	4.7
4	B	1149	ARG	4.4
3	A	1	MET	4.2
4	B	1221	LEU	4.2
4	B	1195	ARG	3.9
4	B	1184	LYS	3.9
4	B	1198	ASN	3.7
4	B	1122	GLN	3.6
4	B	1142	ARG	3.6
4	B	1194	ALA	3.5
4	B	1139	GLU	3.4
4	B	1192	PHE	3.4
4	B	1176	ALA	3.4
4	B	1299	ILE	3.1
4	B	1115	GLU	3.0
4	B	1159	VAL	2.8

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Mol	Chain	Res	Type	RSRZ
3	A	95	LYS	2.8
3	A	59	ILE	2.6
4	B	1177	ASP	2.6
3	A	57	LEU	2.6
4	B	1178	ILE	2.6
4	B	1296	LYS	2.5
4	B	1109	LEU	2.5
3	A	58	LEU	2.5
4	B	1118	ARG	2.5
3	A	41	GLU	2.4
4	B	1124	LYS	2.4
4	B	1120	THR	2.4
4	B	1108	ASN	2.3
4	B	1151	ILE	2.3
4	B	1157	ALA	2.2
3	A	52	ASP	2.2
4	B	1141	VAL	2.1
3	A	148	ILE	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains [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(Å <sup>2</sup> )	Q<0.9
1	5IU	C	1411	20/21	0.97	0.07	11,17,28,42	1
1	5IU	C	1412	20/21	0.97	0.08	15,20,30,55	1

## 6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 6.4 Ligands [i](#)

There are no ligands in this entry.

## 6.5 Other polymers [i](#)

There are no such residues in this entry.