



# Full wwPDB X-ray Structure Validation Report ⓘ

Sep 22, 2021 – 12:23 PM JST

PDB ID : 7CLI  
Title : Structure of NF-kB p52 homodimer bound to P-Selectin kB DNA fragment  
Authors : Meshcheryakov, V.A.; Wang, V.Y.-F.  
Deposited on : 2020-07-21  
Resolution : 3.00 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.

---

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	:	2.23.1
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)	:	2.23.1

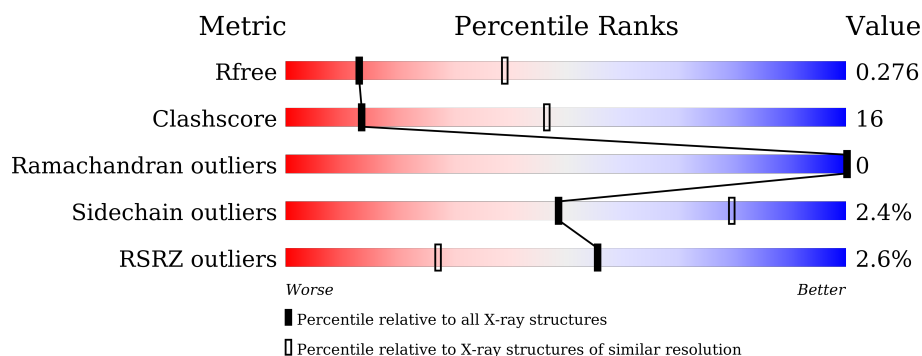
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 3.00 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	2092 (3.00-3.00)
Clashscore	141614	2416 (3.00-3.00)
Ramachandran outliers	138981	2333 (3.00-3.00)
Sidechain outliers	138945	2336 (3.00-3.00)
RSRZ outliers	127900	1990 (3.00-3.00)

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  $\geq 3$ , 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	398	<div> <div>2%</div> <div> <div></div> <div>48%</div> <div>25%</div> <div>27%</div> </div> </div>
1	B	398	<div> <div>2%</div> <div> <div></div> <div>52%</div> <div>22%</div> <div>26%</div> </div> </div>
2	C	18	<div> <div></div> <div> <div>44%</div> <div>50%</div> <div>6%</div> </div> </div>
3	D	18	<div> <div>11%</div> <div> <div></div> <div>72%</div> <div>28%</div> </div> </div>

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 5359 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 Nuclear factor NF-kappa-B p52 subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	292	Total	C	N	O	S	0	0	0
			2300	1449	412	427	12			
1	B	296	Total	C	N	O	S	0	0	0
			2334	1469	420	433	12			

- Molecule 2 is a DNA chain called DNA (5'-D(\*CP\*AP\*AP\*GP\*GP\*GP\*GP\*TP\*CP\*AP\*CP\*CP\*CP\*CP\*TP\*TP\*C)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	17	Total	C	N	O	P	0	0	0
			341	163	62	100	16			

- Molecule 3 is a DNA chain called DNA (5'-D(\*GP\*AP\*AP\*GP\*GP\*GP\*GP\*GP\*TP\*GP\*AP\*CP\*CP\*CP\*CP\*TP\*TP\*G)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	18	Total	C	N	O	P	0	0	0
			372	176	73	106	17			

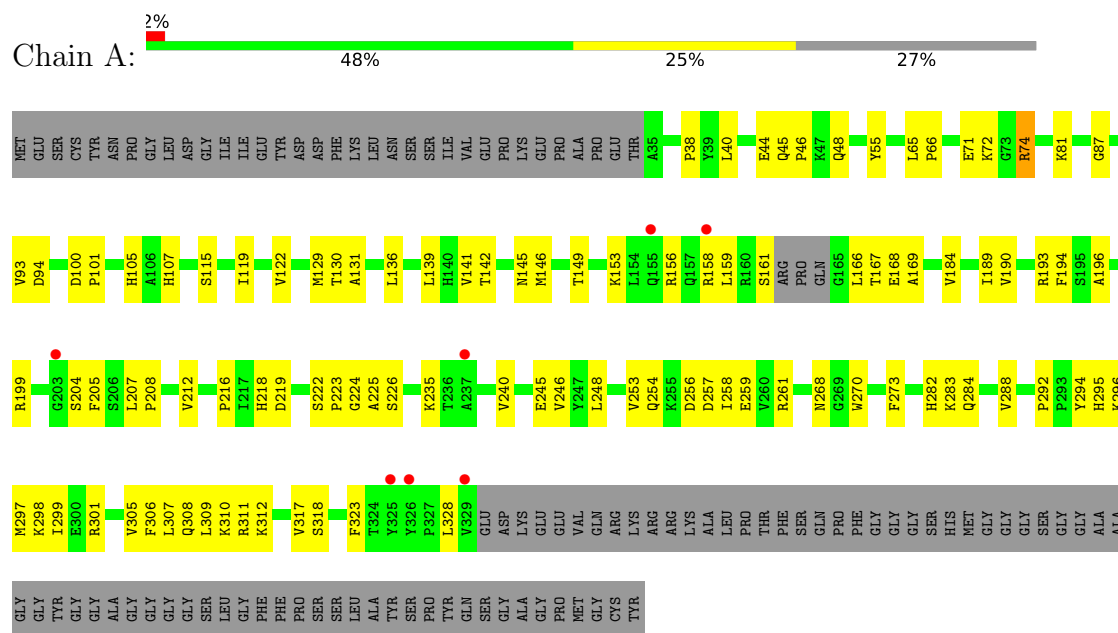
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	6	Total	O	0	0
			6	6		
4	C	1	Total	O	0	0
			1	1		
4	B	4	Total	O	0	0
			4	4		
4	D	1	Total	O	0	0
			1	1		

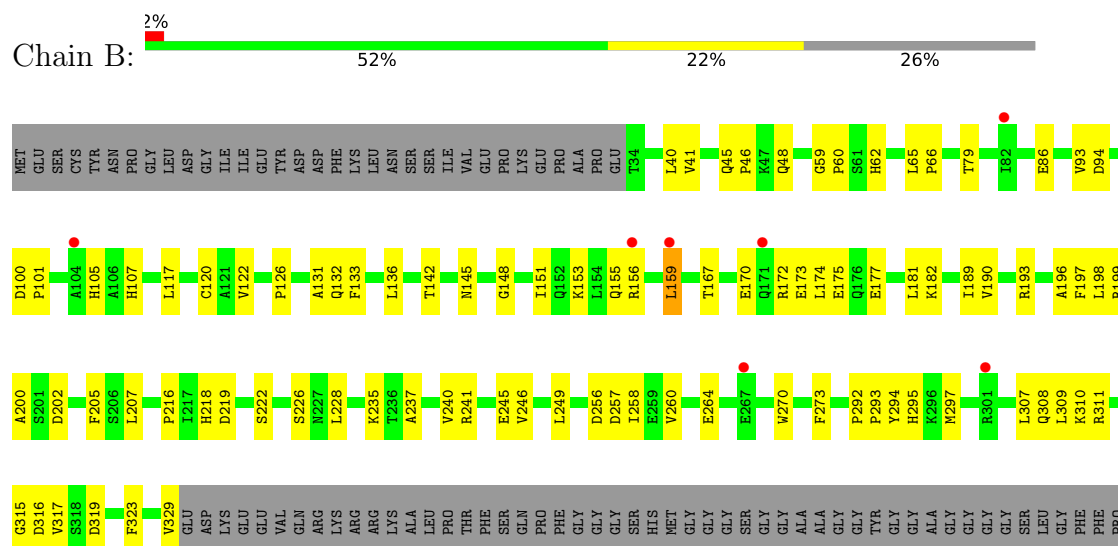
### 3 Residue-property plots

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: Nuclear factor NF-kappa-B p52 subunit



#### • Molecule 1: Nuclear factor NF-kappa-B p52 subunit



SER	SER	LEU	ALA	TYR	SER	PRO	TYR	GLN	SER	GLY	ALA	GLY	PRO	MET	GLY	CYS	TYR
-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----

● Molecule 2: DNA (5'-D(\*CP\*AP\*AP\*GP\*GP\*GP\*GP\*TP\*CP\*AP\*CP\*CP\*CP\*CP\*TP\*TP\*C)-3')



C1	A2	A3	G4	G5	G6	G7	A10	C14	C15	T16	T17	DC
----	----	----	----	----	----	----	-----	-----	-----	-----	-----	----

● Molecule 3: DNA (5'-D(\*GP\*AP\*AP\*GP\*GP\*GP\*GP\*GP\*TP\*GP\*AP\*CP\*CP\*CP\*CP\*TP\*TP\*G)-3')



G1	G4	C14	C15	T16	T17	G18
----	----	-----	-----	-----	-----	-----

## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	84.50Å 85.37Å 140.29Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	42.72 – 3.00 42.69 – 3.00	Depositor EDS
% Data completeness (in resolution range)	94.9 (42.72-3.00) 94.8 (42.69-3.00)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.05 (at 3.01Å)	Xtriage
Refinement program	REFMAC 5.8.0258	Depositor
R, $R_{free}$	0.236 , 0.275 0.238 , 0.276	Depositor DCC
$R_{free}$ test set	973 reflections (4.88%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	86.8	Xtriage
Anisotropy	0.694	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.30 , 73.4	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.000 for k,h,-l	Xtriage
$F_o, F_c$ correlation	0.91	EDS
Total number of atoms	5359	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	115.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 32.31 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 9.5943e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

<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

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	0.71	0/2348	0.92	0/3160
1	B	0.68	0/2384	0.93	0/3211
2	C	0.67	0/381	1.04	0/585
3	D	0.54	0/418	1.02	1/645 (0.2%)
All	All	0.69	0/5531	0.94	1/7601 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	4	DG	C1'-O4'-C4'	-5.81	104.29	110.10

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	2300	0	2309	82	0
1	B	2334	0	2345	74	0
2	C	341	0	192	9	0
3	D	372	0	203	5	0
4	A	6	0	0	0	0
4	B	4	0	0	0	0
4	C	1	0	0	1	0
4	D	1	0	0	0	0

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	5359	0	5049	165	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 (165) 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:270:TRP:NE1	1:A:295:HIS:HB3	1.67	1.10
1:B:167:THR:HG23	1:B:170:GLU:H	1.26	0.97
1:B:240:VAL:HG23	1:B:294:TYR:HB3	1.48	0.93
1:B:308:GLN:NE2	1:B:317:VAL:HG11	1.86	0.91
1:A:270:TRP:CD1	1:A:295:HIS:HB3	2.07	0.88
1:A:158:ARG:HG2	1:A:166:LEU:HB3	1.58	0.85
1:B:155:GLN:O	1:B:159:LEU:HB2	1.78	0.83
1:A:141:VAL:HG11	1:A:149:THR:HG21	1.63	0.80
1:A:158:ARG:HD3	1:A:166:LEU:O	1.83	0.78
1:A:142:THR:HG22	2:C:10:DA:H5'	1.67	0.77
1:A:270:TRP:CG	1:A:295:HIS:ND1	2.52	0.76
1:B:151:ILE:O	1:B:155:GLN:HG3	1.85	0.76
1:B:79:THR:CG2	1:B:132:GLN:HG3	2.14	0.76
1:A:199:ARG:HG3	1:A:204:SER:HB3	1.66	0.76
1:B:79:THR:HG22	1:B:132:GLN:HG3	1.68	0.75
1:B:308:GLN:NE2	1:B:317:VAL:CG1	2.48	0.74
1:B:270:TRP:NE1	1:B:295:HIS:HB3	2.03	0.73
1:A:167:THR:HG22	1:A:169:ALA:H	1.55	0.71
2:C:5:DG:N2	3:D:14:DC:O2	2.20	0.70
1:B:219:ASP:HB3	1:B:222:SER:HB3	1.72	0.69
1:A:189:ILE:HG12	1:A:218:HIS:ND1	2.08	0.69
1:A:283:LYS:O	1:A:284:GLN:HG2	1.94	0.68
1:A:158:ARG:NH1	1:A:168:GLU:HA	2.08	0.67
1:B:167:THR:HG23	1:B:170:GLU:N	2.05	0.67
1:A:199:ARG:HA	1:A:205:PHE:H	1.59	0.67
1:B:235:LYS:HD2	1:B:245:GLU:O	1.95	0.67
1:A:122:VAL:HG21	1:A:131:ALA:HB1	1.74	0.67
1:B:228:LEU:H	1:B:228:LEU:HD12	1.61	0.65
1:A:115:SER:OG	1:A:119:ILE:HG22	1.97	0.65
1:B:122:VAL:HG21	1:B:131:ALA:HB1	1.79	0.64
1:A:40:LEU:HD22	1:A:196:ALA:HB2	1.81	0.63
1:A:270:TRP:CD1	1:A:295:HIS:CB	2.81	0.63
1:B:228:LEU:HD11	1:B:311:ARG:HE	1.65	0.62

*Continued on next page...*



*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:100:ASP:N	1:B:101:PRO:HD2	2.14	0.62
1:A:259:GLU:HG2	1:A:310:LYS:O	2.01	0.60
1:A:297:MET:CE	1:A:328:LEU:HD11	2.31	0.60
1:A:308:GLN:OE1	1:A:317:VAL:HG11	2.01	0.60
1:A:87:GLY:O	1:A:129:MET:HE1	2.01	0.60
1:B:270:TRP:CD1	1:B:295:HIS:HB3	2.36	0.60
1:A:222:SER:HB3	1:A:223:PRO:HD2	1.82	0.60
1:A:161:SER:OG	1:A:166:LEU:HD22	2.02	0.60
1:B:170:GLU:HA	1:B:173:GLU:HB3	1.85	0.59
1:A:254:GLN:HB2	1:A:257:ASP:HB3	1.85	0.58
1:A:297:MET:HE3	1:A:328:LEU:HD11	1.84	0.58
1:B:308:GLN:CD	1:B:317:VAL:CG1	2.73	0.57
1:B:40:LEU:HD23	1:B:41:VAL:N	2.20	0.57
1:A:283:LYS:O	1:A:284:GLN:CG	2.53	0.57
1:A:309:LEU:HB2	1:A:318:SER:HB3	1.86	0.57
1:A:87:GLY:O	1:A:129:MET:CE	2.54	0.56
1:B:170:GLU:HG3	1:B:173:GLU:OE1	2.04	0.56
1:B:235:LYS:HB2	1:B:246:VAL:HG22	1.87	0.56
1:B:200:ALA:HB2	1:B:207:LEU:CD2	2.36	0.56
1:B:79:THR:HG22	1:B:132:GLN:CG	2.35	0.56
1:B:308:GLN:CD	1:B:317:VAL:HG11	2.26	0.56
1:A:55:TYR:CE2	2:C:10:DA:H5''	2.41	0.56
1:B:240:VAL:CG2	1:B:294:TYR:HB3	2.29	0.55
1:A:107:HIS:O	1:A:153:LYS:NZ	2.39	0.55
1:A:161:SER:C	1:A:166:LEU:HD13	2.27	0.55
1:A:48:GLN:HG3	1:A:216:PRO:O	2.06	0.55
1:A:270:TRP:HE1	1:A:295:HIS:HB3	1.69	0.55
1:A:81:LYS:CB	1:A:130:THR:HG22	2.38	0.54
1:B:107:HIS:O	1:B:153:LYS:NZ	2.40	0.54
1:B:240:VAL:HG21	1:B:297:MET:HA	1.88	0.54
1:A:245:GLU:OE1	1:A:288:VAL:HG11	2.07	0.54
1:A:296:LYS:HE3	1:A:298:LYS:HE3	1.91	0.53
2:C:4:DG:O6	1:B:62:HIS:HE1	1.92	0.53
1:A:105:HIS:CD2	1:A:190:VAL:HG12	2.45	0.52
2:C:6:DG:H2'	2:C:7:DG:C8	2.45	0.52
1:A:100:ASP:N	1:A:101:PRO:HD2	2.25	0.52
1:A:270:TRP:CD1	1:A:295:HIS:ND1	2.78	0.51
1:A:161:SER:C	1:A:166:LEU:HD22	2.31	0.51
1:B:189:ILE:HG12	1:B:218:HIS:ND1	2.25	0.51
1:A:81:LYS:HB3	1:A:130:THR:HG22	1.92	0.51
1:B:170:GLU:HG2	1:B:174:LEU:HG	1.91	0.51

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:71:GLU:HG3	1:A:72:LYS:HG2	1.93	0.51
1:A:184:VAL:HG12	1:A:184:VAL:O	2.10	0.51
1:B:86:GLU:HG3	1:B:126:PRO:O	2.11	0.51
1:B:148:GLY:O	1:B:151:ILE:HG13	2.11	0.51
1:B:237:ALA:O	1:B:323:PHE:CZ	2.64	0.51
1:A:45:GLN:OE1	1:A:46:PRO:HD2	2.11	0.50
1:A:296:LYS:O	1:A:299:ILE:HG12	2.12	0.50
1:B:45:GLN:OE1	1:B:46:PRO:HD2	2.12	0.50
1:B:310:LYS:NZ	1:B:315:GLY:O	2.40	0.50
1:B:40:LEU:HD12	1:B:196:ALA:HB2	1.94	0.49
1:B:200:ALA:HB2	1:B:207:LEU:HD21	1.93	0.49
1:A:207:LEU:HD12	1:A:208:PRO:HD2	1.94	0.49
1:A:235:LYS:HB2	1:A:246:VAL:HG22	1.93	0.49
1:B:226:SER:O	1:B:311:ARG:NH2	2.45	0.49
1:B:142:THR:OG1	1:B:145:ASN:HB2	2.13	0.49
1:B:199:ARG:HA	1:B:205:PHE:HA	1.94	0.49
1:A:261:ARG:NH1	1:A:308:GLN:OE1	2.46	0.49
1:B:172:ARG:NH1	1:B:175:GLU:OE2	2.46	0.49
1:A:142:THR:OG1	1:A:145:ASN:HB2	2.13	0.48
1:B:198:LEU:O	1:B:207:LEU:HB2	2.11	0.48
1:B:93:VAL:CG1	1:B:120:CYS:HB2	2.43	0.48
3:D:14:DC:H2'	3:D:15:DC:C5	2.48	0.48
1:B:48:GLN:HG3	1:B:216:PRO:O	2.14	0.48
1:B:105:HIS:CD2	1:B:190:VAL:HG12	2.49	0.48
1:B:167:THR:CG2	1:B:170:GLU:H	2.12	0.47
1:B:273:PHE:O	1:B:292:PRO:HB3	2.15	0.47
1:B:170:GLU:O	1:B:174:LEU:HG	2.15	0.46
1:A:93:VAL:HG22	1:A:194:PHE:CD1	2.50	0.46
1:A:307:LEU:HD23	1:A:323:PHE:HB2	1.97	0.46
1:A:40:LEU:HD22	1:A:196:ALA:CB	2.45	0.46
1:A:74:ARG:O	1:A:74:ARG:HG3	2.15	0.46
1:B:308:GLN:CD	1:B:317:VAL:HG13	2.36	0.46
1:B:94:ASP:OD1	1:B:193:ARG:HB3	2.15	0.46
1:B:199:ARG:HB2	1:B:205:PHE:CD1	2.52	0.46
1:A:40:LEU:HB3	1:A:212:VAL:CG2	2.45	0.45
2:C:5:DG:H2'	2:C:6:DG:O4'	2.16	0.45
1:B:260:VAL:HG22	1:B:309:LEU:HD23	1.98	0.45
1:B:270:TRP:CE2	1:B:295:HIS:HB3	2.52	0.45
1:B:241:ARG:O	1:B:293:PRO:HB3	2.17	0.45
1:B:307:LEU:HD23	1:B:323:PHE:HB2	1.98	0.45
1:A:158:ARG:CD	1:A:166:LEU:O	2.58	0.45

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:79:THR:HG22	1:B:132:GLN:HA	1.99	0.45
1:B:100:ASP:N	1:B:101:PRO:CD	2.78	0.45
1:A:294:TYR:HE2	1:A:299:ILE:HD11	1.82	0.44
1:B:329:VAL:HG13	1:B:329:VAL:O	2.18	0.44
1:A:166:LEU:HD12	1:A:166:LEU:HA	1.90	0.44
1:B:93:VAL:HG12	1:B:120:CYS:HB2	1.99	0.44
1:B:258:ILE:HD11	1:B:309:LEU:HB3	1.99	0.44
1:A:100:ASP:N	1:A:101:PRO:CD	2.81	0.44
1:A:44:GLU:OE2	1:A:44:GLU:HA	2.17	0.44
1:A:305:VAL:HG12	1:A:306:PHE:N	2.33	0.44
1:B:65:LEU:HD12	1:B:66:PRO:HD2	1.99	0.43
1:A:156:ARG:O	1:A:159:LEU:N	2.50	0.43
1:A:224:GLY:O	1:A:311:ARG:NH2	2.51	0.43
1:B:182:LYS:O	1:B:182:LYS:HG2	2.18	0.43
1:A:256:ASP:OD1	1:A:312:LYS:NZ	2.52	0.43
1:B:133:PHE:HB3	1:B:136:LEU:HG	2.01	0.43
1:A:65:LEU:HD12	1:A:66:PRO:HD2	2.01	0.43
1:A:240:VAL:HG23	1:A:297:MET:HE3	2.00	0.43
1:A:273:PHE:O	1:A:292:PRO:HB3	2.18	0.43
2:C:1:DC:H2'	2:C:2:DA:C8	2.54	0.43
1:A:310:LYS:HG3	1:A:317:VAL:HG12	2.01	0.42
1:A:161:SER:OG	1:A:166:LEU:CD2	2.66	0.42
1:A:219:ASP:O	1:A:225:ALA:HB1	2.19	0.42
1:A:253:VAL:HB	1:A:258:ILE:HD13	2.01	0.42
1:B:311:ARG:HD2	1:B:316:ASP:HB2	2.01	0.42
1:B:79:THR:CG2	1:B:132:GLN:CG	2.92	0.42
2:C:14:DC:H5'	4:C:101:HOH:O	2.19	0.42
1:A:283:LYS:C	1:A:284:GLN:HG2	2.39	0.42
1:A:94:ASP:OD1	1:A:193:ARG:HB3	2.20	0.41
1:B:117:LEU:HD22	1:B:156:ARG:HG2	2.01	0.41
1:B:310:LYS:HA	1:B:316:ASP:O	2.19	0.41
3:D:15:DC:H2'	3:D:16:DT:C7	2.50	0.41
1:A:146:MET:O	1:A:149:THR:HG22	2.19	0.41
1:B:240:VAL:HG11	1:B:297:MET:HG3	2.02	0.41
3:D:15:DC:H2''	3:D:16:DT:O5'	2.20	0.41
1:A:139:LEU:C	1:A:139:LEU:HD23	2.40	0.41
1:A:248:LEU:C	1:A:248:LEU:HD23	2.40	0.41
1:A:282:HIS:CE1	1:B:249:LEU:HB3	2.56	0.41
1:B:197:PHE:HA	1:B:207:LEU:O	2.21	0.41
1:A:81:LYS:HB2	1:A:130:THR:HG22	2.02	0.41
2:C:14:DC:H2'	2:C:15:DC:C6	2.56	0.41

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:311:ARG:HD2	1:B:316:ASP:OD2	2.20	0.40
1:A:219:ASP:O	1:A:225:ALA:CB	2.69	0.40
1:A:299:ILE:HG23	1:A:301:ARG:O	2.22	0.40
1:A:305:VAL:N	1:A:323:PHE:O	2.47	0.40
1:A:310:LYS:CG	1:A:317:VAL:HG12	2.52	0.40
1:A:136:LEU:HD23	1:A:136:LEU:HA	1.94	0.40
1:B:59:GLY:HA2	1:B:60:PRO:HD3	1.87	0.40
1:B:189:ILE:CD1	1:B:218:HIS:CE1	3.05	0.40
3:D:16:DT:H2'	3:D:17:DT:C6	2.56	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	288/398 (72%)	275 (96%)	13 (4%)	0	100	100
1	B	294/398 (74%)	284 (97%)	10 (3%)	0	100	100
All	All	582/796 (73%)	559 (96%)	23 (4%)	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	253/330 (77%)	249 (98%)	4 (2%)	62	86
1	B	257/330 (78%)	249 (97%)	8 (3%)	40	75
All	All	510/660 (77%)	498 (98%)	12 (2%)	49	79

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	38	PRO
1	A	74	ARG
1	A	226	SER
1	A	268	ASN
1	B	159	LEU
1	B	177	GLU
1	B	181	LEU
1	B	202	ASP
1	B	256	ASP
1	B	257	ASP
1	B	264	GLU
1	B	319	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	227	ASN
1	B	295	HIS

### 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 [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 [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	292/398 (73%)	0.14	7 (2%) 59 30	73, 109, 167, 218	0
1	B	296/398 (74%)	0.14	7 (2%) 59 30	66, 110, 175, 212	0
2	C	17/18 (94%)	0.29	0 100 100	89, 110, 146, 153	0
3	D	18/18 (100%)	1.04	2 (11%) 5 1	88, 112, 157, 157	0
All	All	623/832 (74%)	0.17	16 (2%) 56 27	66, 110, 171, 218	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	D	16	DT	3.5
3	D	1	DG	3.5
1	A	155	GLN	3.4
1	A	158	ARG	3.3
1	A	203	GLY	3.1
1	B	301	ARG	3.1
1	B	159	LEU	3.0
1	A	325	TYR	2.7
1	B	267	GLU	2.7
1	B	82	ILE	2.4
1	A	329	VAL	2.4
1	A	326	TYR	2.3
1	B	104	ALA	2.3
1	B	171	GLN	2.2
1	B	156	ARG	2.2
1	A	237	ALA	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 monosaccharides 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.