



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2018 – 11:40 am GMT

PDB ID : 1D0E  
Title : CRYSTAL STRUCTURES OF THE N-TERMINAL FRAGMENT FROM  
MOLONEY MURINE LEUKEMIA VIRUS REVERSE TRANSCRIPTASE  
COMPLEXED WITH NUCLEIC ACID: FUNCTIONAL IMPLICATIONS  
FOR TEMPLATE-PRIMER BINDING TO THE FINGERS DOMAIN  
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Deposited on : 1999-09-09  
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.

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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	:	trunk30967
Percentile statistics	:	20171227.v01 (using entries in the PDB archive December 27th 2017)
Refmac	:	5.8.0158
CCP4	:	7.0 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk30967

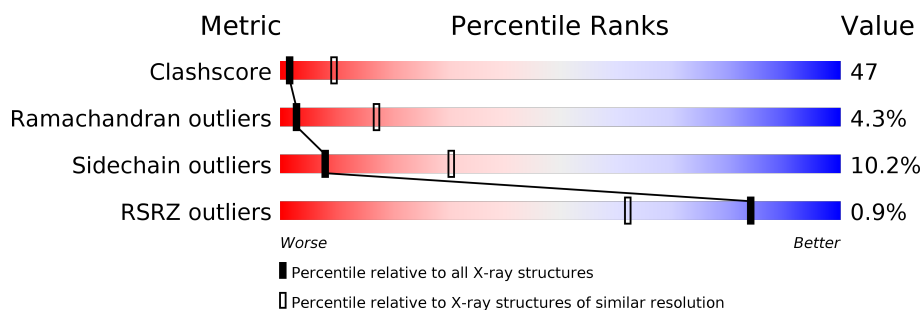
# 1 Overall quality at a glance

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))
Clashscore	122126	2167 (3.00-3.00)
Ramachandran outliers	120053	2101 (3.00-3.00)
Sidechain outliers	120020	2104 (3.00-3.00)
RSRZ outliers	108989	1751 (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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	E	11	
1	F	11	
2	A	259	
2	B	259	

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4570 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(\*TP\*TP\*TP\*CP\*AP\*TP\*GP\*CP\*AP\*TP\*G)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	E	11	Total	C	N	O	P	0	0	0
			221	108	36	67	10			
1	F	11	Total	C	N	O	P	0	0	0
			221	108	36	67	10			

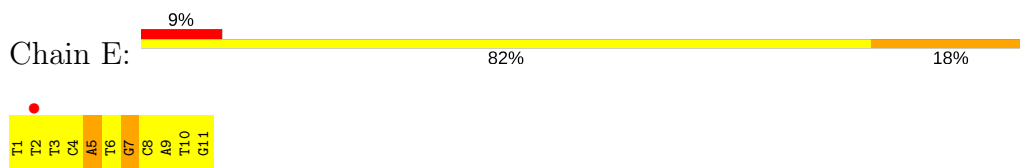
- Molecule 2 is a protein called REVERSE TRANSCRIPTASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	A	259	Total	C	N	O	S	0	0	0
			2069	1327	362	372	8			
2	B	257	Total	C	N	O	S	0	0	0
			2059	1322	360	369	8			

### 3 Residue-property plots

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.

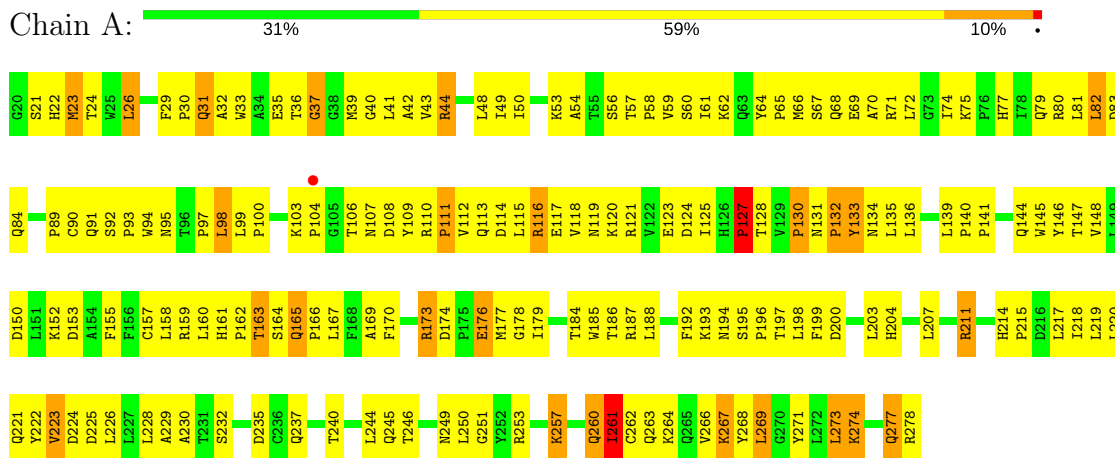
- Molecule 1: 5'-D(\*TP\*TP\*TP\*CP\*AP\*TP\*GP\*CP\*AP\*TP\*G)-3'



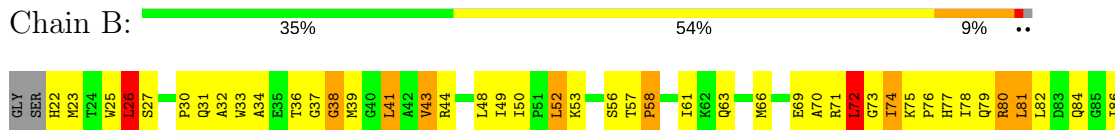
- Molecule 1: 5'-D(\*TP\*TP\*TP\*CP\*AP\*TP\*GP\*CP\*AP\*TP\*G)-3'



- Molecule 2: REVERSE TRANSCRIPTASE



- Molecule 2: REVERSE TRANSCRIPTASE





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	66.16Å 63.54Å 73.50Å 90.00° 103.85° 90.00°	Depositor
Resolution (Å)	500.00 – 3.00 28.48 – 3.00	Depositor EDS
% Data completeness (in resolution range)	95.7 (500.00-3.00) 95.9 (28.48-3.00)	Depositor EDS
$R_{merge}$	0.15	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.39 (at 3.00Å)	Xtriage
Refinement program	CNS	Depositor
R, $R_{free}$	0.215 , 0.298 0.210 , (Not available)	Depositor DCC
$R_{free}$ test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	31.2	Xtriage
Anisotropy	0.747	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 57.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.43$ , $\langle L^2 \rangle = 0.26$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.89	EDS
Total number of atoms	4570	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.90% 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

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	E	0.51	0/246	0.68	0/378
1	F	0.44	0/246	0.76	0/378
2	A	0.63	5/2126 (0.2%)	0.89	7/2896 (0.2%)
2	B	0.56	1/2116 (0.0%)	0.95	5/2883 (0.2%)
All	All	0.59	6/4734 (0.1%)	0.90	12/6535 (0.2%)

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	E	0	2

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	273	LEU	CA-C	9.15	1.76	1.52
2	A	274	LYS	N-CA	7.89	1.62	1.46
2	A	127	PRO	N-CA	7.48	1.59	1.47
2	A	261	ILE	C-N	-6.50	1.19	1.34
2	B	107	ASN	N-CA	5.96	1.58	1.46
2	A	273	LEU	C-N	5.45	1.46	1.34

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	27	SER	N-CA-CB	-21.60	78.10	110.50
2	A	127	PRO	CA-N-CD	-11.83	94.94	111.50
2	B	26	LEU	CB-CA-C	11.12	131.33	110.20
2	A	273	LEU	CA-C-N	7.62	133.96	117.20
2	A	127	PRO	N-CA-C	7.21	130.84	112.10
2	A	273	LEU	N-CA-C	6.99	129.87	111.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	260	GLN	C-N-CA	6.09	136.93	121.70
2	B	27	SER	N-CA-C	5.55	125.97	111.00
2	B	99	LEU	C-N-CD	-5.54	108.42	120.60
2	A	260	GLN	O-C-N	-5.34	114.16	122.70
2	B	72	LEU	CA-CB-CG	5.25	127.39	115.30
2	A	127	PRO	N-CD-CG	5.04	110.76	103.20

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	E	5	DA	Sidechain
1	E	7	DG	Sidechain

## 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	E	221	0	128	20	0
1	F	221	0	128	11	0
2	A	2069	0	2079	204	0
2	B	2059	0	2072	194	0
All	All	4570	0	4407	416	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 47.

All (416) 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:A:273:LEU:CA	2:A:273:LEU:C	1.76	1.52
2:A:264:LYS:O	2:A:274:LYS:HB2	1.39	1.20
2:A:68:GLN:HE21	2:A:69:GLU:HG3	1.18	1.08
2:B:74:ILE:HG13	2:B:111:PRO:HG3	1.40	1.01
2:A:266:VAL:HG22	2:A:273:LEU:HD21	1.44	0.97

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:165:GLN:HE21	2:A:186:THR:HA	1.25	0.97
2:B:131:ASN:HB3	2:B:134:ASN:HB3	1.44	0.97
2:B:201:GLU:HG2	2:B:205:ARG:HH22	1.30	0.95
1:E:5:DA:H2''	1:E:6:DT:H5'	1.50	0.93
2:B:74:ILE:CD1	2:B:111:PRO:HD3	2.00	0.92
1:E:6:DT:H3	1:F:9:DA:H61	1.12	0.91
2:B:26:LEU:HD11	2:B:36:THR:HG23	1.52	0.90
2:B:74:ILE:HD11	2:B:111:PRO:HD3	1.54	0.88
2:A:232:SER:HB3	2:A:235:ASP:OD2	1.72	0.88
1:E:9:DA:H4'	1:E:10:DT:OP1	1.74	0.86
2:A:214:HIS:HB3	2:A:217:LEU:HD12	1.57	0.86
2:A:68:GLN:NE2	2:A:69:GLU:HG3	1.91	0.86
2:B:61:ILE:HD11	2:B:117:GLU:HG3	1.57	0.86
2:A:266:VAL:CG2	2:A:273:LEU:HD21	2.05	0.86
2:A:266:VAL:O	2:A:273:LEU:HD22	1.76	0.85
2:B:71:ARG:HH22	2:B:75:LYS:HZ3	1.24	0.84
2:B:101:VAL:O	2:B:109:TYR:CD1	2.31	0.83
2:B:101:VAL:O	2:B:109:TYR:HD1	1.61	0.83
2:B:71:ARG:HH22	2:B:75:LYS:NZ	1.77	0.81
2:A:31:GLN:HE21	2:A:31:GLN:HA	1.46	0.81
2:B:149:LEU:HD12	2:B:149:LEU:O	1.79	0.81
2:A:113:GLN:HE21	2:A:170:PHE:HB3	1.46	0.81
2:A:91:GLN:HG3	2:A:91:GLN:O	1.80	0.81
2:A:21:SER:OG	2:A:24:THR:HG23	1.81	0.80
2:A:266:VAL:H	2:A:273:LEU:CD2	1.93	0.80
2:B:43:VAL:CG2	2:B:44:ARG:HH11	1.94	0.80
2:B:74:ILE:CG1	2:B:111:PRO:HG3	2.13	0.79
2:A:266:VAL:H	2:A:273:LEU:HD23	1.49	0.77
2:B:225:ASP:O	2:B:226:LEU:HD23	1.85	0.77
2:A:48:LEU:HD13	2:A:198:LEU:HD13	1.66	0.77
2:B:118:VAL:O	2:B:122:VAL:HG23	1.84	0.77
1:F:9:DA:H8	1:F:9:DA:OP2	1.68	0.77
2:A:77:HIS:O	2:A:81:LEU:HD23	1.84	0.76
2:B:266:VAL:O	2:B:272:LEU:HD12	1.85	0.76
2:A:41:LEU:HD13	2:B:43:VAL:HB	1.65	0.76
2:B:272:LEU:O	2:B:273:LEU:HD23	1.86	0.76
1:F:8:DC:H2''	1:F:9:DA:OP2	1.84	0.76
2:A:146:TYR:CE2	2:A:273:LEU:HG	2.20	0.76
2:B:71:ARG:NH2	2:B:75:LYS:HZ3	1.84	0.75
2:A:130:PRO:HB2	2:A:135:LEU:HD22	1.69	0.75
2:A:39:MET:SD	2:A:84:GLN:HG3	2.26	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:70:ALA:O	2:A:74:ILE:HG12	1.85	0.75
2:A:165:GLN:NE2	2:A:186:THR:HA	2.02	0.73
2:A:266:VAL:O	2:A:273:LEU:CD2	2.37	0.73
1:E:6:DT:H2"	1:E:7:DG:N7	2.04	0.73
2:A:203:LEU:HD11	2:A:207:LEU:HD11	1.71	0.72
2:B:43:VAL:HG21	2:B:44:ARG:HH11	1.55	0.72
2:A:155:PHE:HB3	2:A:188:LEU:HB3	1.70	0.72
2:B:26:LEU:CD1	2:B:36:THR:HG23	2.19	0.72
2:A:130:PRO:O	2:A:131:ASN:C	2.24	0.71
2:B:159:ARG:HA	2:B:187:ARG:HA	1.72	0.71
2:A:127:PRO:O	2:A:128:THR:OG1	2.09	0.71
2:B:38:GLY:O	2:B:253:ARG:HG3	1.90	0.71
2:B:102:LYS:HD3	2:B:102:LYS:O	1.91	0.70
2:A:41:LEU:HD21	2:A:159:ARG:HD2	1.72	0.70
2:B:218:ILE:HB	2:B:229:ALA:HB3	1.72	0.70
2:B:74:ILE:HD11	2:B:111:PRO:CD	2.22	0.69
2:A:152:LYS:HG3	2:A:153:ASP:OD1	1.92	0.69
2:B:90:CYS:SG	2:B:184:THR:HB	2.31	0.69
2:B:214:HIS:HB3	2:B:217:LEU:HD12	1.73	0.69
2:A:161:HIS:CD2	2:A:163:THR:OG1	2.45	0.69
2:B:73:GLY:O	2:B:76:PRO:HD2	1.93	0.69
2:B:61:ILE:CD1	2:B:117:GLU:HG3	2.22	0.69
2:A:273:LEU:CB	2:A:273:LEU:C	2.63	0.67
2:A:214:HIS:CB	2:A:217:LEU:HD12	2.24	0.66
2:B:113:GLN:NE2	2:B:170:PHE:HB3	2.10	0.66
2:A:26:LEU:HD22	2:A:33:TRP:CD1	2.31	0.66
2:A:92:SER:CB	2:A:166:PRO:HB3	2.25	0.66
2:A:90:CYS:SG	2:A:184:THR:HB	2.36	0.66
2:A:114:ASP:CG	2:A:116:ARG:HE	1.99	0.66
2:B:39:MET:CE	2:B:86:ILE:HD11	2.26	0.66
2:B:113:GLN:HE21	2:B:170:PHE:HB3	1.61	0.65
2:B:34:ALA:HA	2:B:253:ARG:HB3	1.78	0.65
2:B:98:LEU:O	2:B:100:PRO:HD3	1.96	0.65
2:B:206:ASP:O	2:B:208:ALA:N	2.28	0.65
2:B:90:CYS:O	2:B:182:GLN:HG3	1.97	0.65
2:A:71:ARG:HH22	2:A:173:ARG:H	1.45	0.65
2:A:162:PRO:C	2:A:164:SER:H	2.01	0.64
2:A:146:TYR:CG	2:A:273:LEU:HD11	2.33	0.64
2:A:203:LEU:O	2:A:207:LEU:HG	1.98	0.64
2:A:80:ARG:HG3	2:A:81:LEU:HD22	1.79	0.64
2:B:156:PHE:HA	2:B:187:ARG:HB3	1.80	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:53:LYS:HE3	2:B:122:VAL:O	1.98	0.64
2:B:195:SER:HB2	2:B:196:PRO:HD3	1.80	0.64
2:A:83:ASP:OD1	2:B:161:HIS:HE1	1.80	0.64
2:A:161:HIS:HD2	2:A:163:THR:OG1	1.80	0.64
2:A:148:VAL:HB	2:A:260:GLN:HB2	1.80	0.64
2:B:201:GLU:HG2	2:B:205:ARG:NH2	2.09	0.64
2:A:273:LEU:C	2:A:273:LEU:HA	2.08	0.63
2:B:205:ARG:HG3	2:B:205:ARG:HH11	1.63	0.63
1:E:2:DT:H3	1:E:4:DC:H41	1.46	0.63
1:F:4:DC:H2''	1:F:5:DA:H5'	1.81	0.63
1:F:9:DA:OP2	1:F:9:DA:C8	2.50	0.63
2:A:113:GLN:NE2	2:A:170:PHE:HB3	2.15	0.62
2:B:161:HIS:HD2	2:B:163:THR:HB	1.64	0.62
2:A:164:SER:O	2:A:167:LEU:HG	1.99	0.62
1:E:5:DA:C2'	1:E:6:DT:H5'	2.27	0.62
2:B:22:HIS:HD2	2:B:23:MET:H	1.48	0.62
2:B:56:SER:OG	2:B:121:ARG:HB3	1.99	0.61
2:B:133:TYR:O	2:B:135:LEU:N	2.34	0.61
1:E:6:DT:H3	1:F:9:DA:N6	1.92	0.61
2:A:147:THR:HA	2:A:260:GLN:O	2.00	0.61
2:A:43:VAL:HG22	2:A:250:LEU:O	2.00	0.61
2:B:33:TRP:CD2	2:B:256:ALA:HB2	2.36	0.61
2:B:165:GLN:HB2	2:B:166:PRO:HD3	1.83	0.60
2:A:121:ARG:HH11	2:A:121:ARG:HG3	1.67	0.60
2:A:22:HIS:HB3	2:A:23:MET:CE	2.31	0.60
2:A:106:THR:C	2:A:108:ASP:H	2.05	0.60
2:A:125:ILE:O	2:A:127:PRO:HD3	2.02	0.60
2:B:101:VAL:O	2:B:109:TYR:HA	2.00	0.60
2:B:74:ILE:HA	2:B:77:HIS:ND1	2.15	0.60
2:A:74:ILE:HD13	2:A:100:PRO:HB3	1.82	0.60
2:B:61:ILE:HD11	2:B:117:GLU:CG	2.31	0.60
2:A:115:LEU:O	2:A:119:ASN:ND2	2.35	0.59
2:B:50:ILE:HG12	2:B:164:SER:HB3	1.84	0.59
2:B:162:PRO:HA	2:B:165:GLN:HG3	1.84	0.59
2:A:266:VAL:CG2	2:A:273:LEU:CD2	2.80	0.59
2:A:26:LEU:HA	2:A:33:TRP:CD1	2.37	0.59
2:A:92:SER:HB3	2:A:166:PRO:HB3	1.83	0.59
2:B:41:LEU:O	2:B:251:GLY:HA3	2.03	0.59
2:A:114:ASP:OD2	2:A:116:ARG:NE	2.32	0.58
2:B:26:LEU:CD1	2:B:36:THR:CG2	2.79	0.58
2:A:49:ILE:O	2:A:50:ILE:HD13	2.03	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:272:LEU:HD12	2:B:273:LEU:H	1.68	0.58
2:B:72:LEU:HD13	2:B:73:GLY:H	1.67	0.58
2:B:221:GLN:HB2	2:B:226:LEU:HD22	1.86	0.58
2:A:32:ALA:O	2:A:253:ARG:HA	2.03	0.58
2:A:79:GLN:NE2	2:A:177:MET:SD	2.77	0.58
2:A:136:LEU:HD12	2:A:136:LEU:N	2.18	0.58
2:B:175:PRO:O	2:B:176:GLU:C	2.42	0.58
2:B:115:LEU:O	2:B:119:ASN:ND2	2.37	0.57
2:A:145:TRP:O	2:A:229:ALA:HA	2.04	0.57
2:B:151:LEU:HD23	2:B:254:ALA:HB2	1.85	0.57
2:A:89:PRO:HG3	2:A:179:ILE:HD12	1.87	0.57
1:E:1:DT:O5'	2:A:104:PRO:HG3	2.04	0.57
2:A:132:PRO:O	2:A:133:TYR:C	2.42	0.56
2:A:71:ARG:O	2:A:74:ILE:HG13	2.04	0.56
2:A:200:ASP:HB2	2:A:223:VAL:HG22	1.86	0.56
2:A:58:PRO:HG3	2:A:94:TRP:HE1	1.69	0.56
2:B:221:GLN:HB2	2:B:226:LEU:CD2	2.34	0.56
2:A:266:VAL:N	2:A:273:LEU:CD2	2.66	0.56
2:B:39:MET:HE3	2:B:86:ILE:HD11	1.87	0.56
2:A:174:ASP:OD1	2:A:176:GLU:HB3	2.06	0.56
2:A:98:LEU:HD23	2:A:170:PHE:HB2	1.87	0.56
2:B:22:HIS:CD2	2:B:23:MET:N	2.74	0.55
2:A:91:GLN:HG2	2:B:89:PRO:HG2	1.88	0.55
2:A:22:HIS:HB3	2:A:23:MET:HE3	1.87	0.55
2:B:204:HIS:N	2:B:221:GLN:HE22	2.04	0.55
2:A:165:GLN:HB3	2:A:185:TRP:O	2.07	0.55
2:A:29:PHE:N	2:A:30:PRO:HD3	2.21	0.55
2:A:49:ILE:N	2:A:49:ILE:HD12	2.21	0.55
2:A:158:LEU:O	2:A:187:ARG:HA	2.07	0.55
2:A:273:LEU:O	2:A:273:LEU:CA	2.48	0.55
2:A:64:TYR:CE2	2:A:99:LEU:HD21	2.42	0.55
2:B:75:LYS:HZ2	2:B:176:GLU:HG2	1.71	0.55
2:B:175:PRO:O	2:B:177:MET:N	2.40	0.55
2:A:120:LYS:HZ3	2:A:121:ARG:HH12	1.53	0.54
2:A:64:TYR:HE2	2:A:99:LEU:HD21	1.73	0.54
2:A:26:LEU:HD11	2:A:36:THR:OG1	2.08	0.54
2:A:192:PHE:HB3	2:A:195:SER:OG	2.07	0.54
2:A:261:ILE:HD12	2:A:261:ILE:N	2.22	0.54
1:E:8:DC:H2''	1:E:9:DA:C8	2.43	0.54
2:A:119:ASN:O	2:A:193:LYS:NZ	2.32	0.54
2:A:273:LEU:HD23	2:A:273:LEU:H	1.72	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:125:ILE:HG12	2:B:194:ASN:OD1	2.08	0.54
2:B:197:THR:O	2:B:201:GLU:HB2	2.08	0.54
2:B:78:ILE:HG23	2:B:87:LEU:CD1	2.37	0.54
2:A:117:GLU:HA	2:A:117:GLU:OE1	2.08	0.54
2:A:157:CYS:C	2:A:158:LEU:HD12	2.28	0.54
2:B:102:LYS:C	2:B:102:LYS:HD3	2.28	0.54
2:A:127:PRO:C	2:A:128:THR:OG1	2.45	0.54
2:A:71:ARG:NH2	2:A:173:ARG:H	2.05	0.54
2:B:258:LYS:NZ	2:B:258:LYS:HB3	2.23	0.54
2:B:72:LEU:HD13	2:B:73:GLY:N	2.22	0.54
2:B:139:LEU:HD13	2:B:220:LEU:HD21	1.90	0.53
2:A:226:LEU:N	2:A:226:LEU:HD12	2.23	0.53
2:A:31:GLN:CA	2:A:31:GLN:HE21	2.19	0.53
2:B:66:MET:SD	2:B:100:PRO:HG3	2.49	0.53
1:F:11:DG:O3'	2:A:115:LEU:N	2.42	0.53
2:B:179:ILE:HD13	2:B:183:LEU:HD21	1.91	0.53
2:B:92:SER:N	2:B:182:GLN:OE1	2.41	0.53
2:B:74:ILE:HD12	2:B:111:PRO:HD3	1.87	0.53
2:A:200:ASP:HB2	2:A:223:VAL:HA	1.91	0.52
2:B:154:ALA:HB1	2:B:199:PHE:CZ	2.44	0.52
2:B:74:ILE:HG13	2:B:111:PRO:CG	2.27	0.52
1:E:1:DT:H3'	1:E:2:DT:C5'	2.40	0.52
2:A:42:ALA:O	2:A:159:ARG:NH1	2.42	0.52
2:A:266:VAL:N	2:A:273:LEU:HD21	2.25	0.52
1:F:3:DT:H2''	1:F:4:DC:OP2	2.09	0.52
2:A:92:SER:HB2	2:A:166:PRO:HB3	1.91	0.52
2:B:168:PHE:HB3	2:B:185:TRP:HD1	1.75	0.52
2:B:22:HIS:HD2	2:B:23:MET:N	2.08	0.52
2:A:266:VAL:H	2:A:273:LEU:HD21	1.74	0.52
2:A:146:TYR:HA	2:A:228:LEU:O	2.10	0.52
2:B:37:GLY:O	2:B:38:GLY:O	2.27	0.52
2:A:26:LEU:HA	2:A:33:TRP:NE1	2.25	0.51
2:B:103:LYS:HE3	2:B:104:PRO:HD2	1.90	0.51
2:A:193:LYS:HG3	2:A:194:ASN:N	2.25	0.51
2:B:66:MET:HB3	2:B:70:ALA:HB3	1.92	0.51
2:A:83:ASP:OD1	2:B:161:HIS:CE1	2.61	0.51
2:B:258:LYS:HB3	2:B:258:LYS:HZ3	1.74	0.51
2:B:30:PRO:HA	2:B:36:THR:OG1	2.11	0.51
2:B:71:ARG:HH12	2:B:75:LYS:HE2	1.74	0.51
2:A:269:LEU:C	2:A:271:TYR:H	2.14	0.51
2:A:53:LYS:HD2	2:A:121:ARG:HA	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:209:ASP:HA	2:B:212:ILE:CD1	2.40	0.51
2:A:173:ARG:NH1	2:A:173:ARG:HB2	2.26	0.51
2:B:176:GLU:OE1	2:B:176:GLU:N	2.44	0.51
2:B:215:PRO:HG2	2:B:216:ASP:OD1	2.10	0.51
2:A:58:PRO:HG3	2:A:94:TRP:NE1	2.26	0.51
2:B:49:ILE:HD12	2:B:49:ILE:N	2.26	0.51
2:B:39:MET:HE1	2:B:86:ILE:HD11	1.92	0.51
2:A:120:LYS:O	2:A:120:LYS:HG3	2.11	0.51
2:A:277:GLN:CD	2:A:278:ARG:H	2.14	0.51
2:B:75:LYS:N	2:B:76:PRO:HD2	2.26	0.51
2:B:249:ASN:N	2:B:249:ASN:ND2	2.58	0.50
2:A:44:ARG:HA	2:B:31:GLN:NE2	2.26	0.50
2:A:223:VAL:HG12	2:A:224:ASP:N	2.26	0.50
2:A:74:ILE:HG23	2:A:109:TYR:HB3	1.93	0.50
2:A:261:ILE:O	2:A:262:CYS:C	2.47	0.50
2:B:133:TYR:C	2:B:135:LEU:H	2.15	0.50
2:B:41:LEU:HD21	2:B:159:ARG:NH1	2.26	0.50
2:B:268:TYR:CE2	2:B:269:LEU:HD22	2.47	0.50
2:A:136:LEU:H	2:A:136:LEU:CD1	2.24	0.50
2:A:35:GLU:OE2	2:A:257:LYS:HD3	2.11	0.50
2:A:72:LEU:HA	2:A:75:LYS:HB2	1.93	0.50
2:B:204:HIS:CE1	2:B:221:GLN:HG2	2.47	0.50
2:B:80:ARG:O	2:B:84:GLN:HG3	2.12	0.50
2:A:56:SER:HB3	2:A:121:ARG:HB3	1.94	0.50
2:B:175:PRO:C	2:B:177:MET:N	2.65	0.50
2:B:57:THR:HG23	2:B:58:PRO:HD2	1.93	0.50
2:A:221:GLN:HG3	2:A:221:GLN:O	2.11	0.50
2:B:66:MET:CE	2:B:100:PRO:HD3	2.42	0.50
2:A:128:THR:OG1	2:A:197:THR:HG23	2.12	0.49
1:F:4:DC:H2''	1:F:5:DA:C5'	2.42	0.49
2:A:95:ASN:HA	2:A:169:ALA:O	2.13	0.49
2:A:68:GLN:O	2:A:72:LEU:HG	2.12	0.49
2:A:98:LEU:HD23	2:A:170:PHE:CB	2.42	0.49
2:A:150:ASP:OD2	2:A:152:LYS:NZ	2.46	0.49
2:A:178:GLY:C	2:A:179:ILE:HD13	2.33	0.49
2:B:209:ASP:HA	2:B:212:ILE:HD12	1.95	0.49
2:B:106:THR:O	2:B:106:THR:HG23	2.13	0.48
2:B:227:LEU:HB2	2:B:268:TYR:CE1	2.48	0.48
2:A:195:SER:HB2	2:A:196:PRO:HD3	1.95	0.48
2:B:143:HIS:CE1	2:B:231:THR:HG23	2.47	0.48
2:A:35:GLU:OE1	2:A:35:GLU:N	2.41	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:172:TRP:CE3	2:B:172:TRP:HA	2.48	0.48
2:B:206:ASP:C	2:B:208:ALA:H	2.16	0.48
1:E:4:DC:O5'	1:E:4:DC:H6	1.97	0.48
2:A:58:PRO:HA	2:A:94:TRP:CE2	2.49	0.48
2:B:71:ARG:HH12	2:B:75:LYS:CE	2.27	0.48
2:A:61:ILE:O	2:A:95:ASN:ND2	2.40	0.48
2:B:53:LYS:NZ	2:B:120:LYS:O	2.39	0.47
2:A:240:THR:O	2:A:244:LEU:HG	2.14	0.47
2:B:205:ARG:CG	2:B:205:ARG:HH11	2.27	0.47
2:A:162:PRO:O	2:A:164:SER:N	2.47	0.47
2:A:140:PRO:HD2	2:A:218:ILE:CD1	2.45	0.47
2:B:74:ILE:O	2:B:77:HIS:HB2	2.14	0.47
2:A:53:LYS:HB2	2:A:56:SER:OG	2.15	0.47
2:B:71:ARG:NH1	2:B:75:LYS:HE2	2.29	0.47
2:A:123:GLU:HA	2:A:123:GLU:OE2	2.14	0.47
1:E:5:DA:H2''	1:E:6:DT:C5'	2.35	0.47
2:A:136:LEU:N	2:A:136:LEU:CD1	2.78	0.47
2:A:199:PHE:HE2	2:A:223:VAL:O	1.98	0.47
2:A:97:PRO:HD2	2:A:114:ASP:HB3	1.97	0.47
2:B:114:ASP:O	2:B:114:ASP:OD1	2.32	0.47
2:B:116:ARG:HA	2:B:119:ASN:HD22	1.80	0.47
2:B:98:LEU:O	2:B:100:PRO:CD	2.61	0.47
2:B:32:ALA:O	2:B:253:ARG:HA	2.15	0.47
2:A:41:LEU:N	2:A:251:GLY:O	2.44	0.46
2:B:169:ALA:HA	2:B:183:LEU:O	2.15	0.46
2:B:143:HIS:ND1	2:B:231:THR:HG23	2.30	0.46
2:B:249:ASN:N	2:B:249:ASN:HD22	2.13	0.46
1:E:6:DT:O2	1:F:9:DA:N1	2.48	0.46
2:B:96:THR:OG1	2:B:113:GLN:NE2	2.48	0.46
2:A:192:PHE:CE2	2:A:194:ASN:HB2	2.51	0.46
2:A:173:ARG:CB	2:A:173:ARG:CZ	2.94	0.46
2:B:71:ARG:NH2	2:B:75:LYS:NZ	2.50	0.46
2:B:133:TYR:C	2:B:135:LEU:N	2.69	0.46
2:B:78:ILE:HG22	2:B:82:LEU:CD2	2.46	0.46
1:E:9:DA:H1'	1:E:10:DT:H5'	1.96	0.46
2:A:98:LEU:HD23	2:A:170:PHE:CG	2.51	0.46
2:B:265:GLN:HG3	2:B:274:LYS:HG3	1.98	0.46
2:A:173:ARG:HB3	2:A:173:ARG:CZ	2.45	0.46
2:A:211:ARG:HD3	2:A:219:LEU:HB3	1.96	0.46
2:A:267:LYS:O	2:A:267:LYS:HE3	2.15	0.46
2:B:172:TRP:HE3	2:B:172:TRP:HA	1.79	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:188:LEU:HD21	2:B:198:LEU:HB2	1.97	0.46
1:F:3:DT:O2	1:F:3:DT:O4'	2.34	0.46
2:B:205:ARG:HB2	2:B:205:ARG:CZ	2.46	0.46
2:B:73:GLY:O	2:B:76:PRO:CD	2.64	0.45
2:A:66:MET:HE3	2:A:71:ARG:HH11	1.81	0.45
2:B:33:TRP:CZ3	2:B:256:ALA:HA	2.52	0.45
2:B:227:LEU:HB2	2:B:268:TYR:CD1	2.52	0.45
2:B:71:ARG:HH12	2:B:75:LYS:NZ	2.15	0.45
2:A:218:ILE:HG22	2:A:220:LEU:CD1	2.47	0.45
2:A:266:VAL:HG23	2:A:273:LEU:CD2	2.46	0.45
2:A:59:VAL:O	2:A:59:VAL:HG13	2.17	0.45
2:B:146:TYR:HB2	2:B:266:VAL:HG11	1.98	0.45
2:B:154:ALA:O	2:B:157:CYS:HB2	2.17	0.45
2:B:167:LEU:HD13	2:B:167:LEU:H	1.81	0.45
2:B:154:ALA:HB1	2:B:199:PHE:CE2	2.52	0.45
2:B:89:PRO:O	2:B:90:CYS:HB3	2.17	0.45
2:A:141:PRO:O	2:A:144:GLN:HG3	2.17	0.45
2:A:246:THR:HG22	2:A:250:LEU:HD12	1.98	0.45
1:E:3:DT:H4'	1:E:4:DC:OP1	2.15	0.45
2:A:222:TYR:O	2:A:223:VAL:HB	2.17	0.44
2:B:101:VAL:O	2:B:109:TYR:CA	2.65	0.44
2:B:71:ARG:O	2:B:72:LEU:C	2.56	0.44
2:A:245:GLN:NE2	2:A:249:ASN:OD1	2.43	0.44
2:B:78:ILE:O	2:B:81:LEU:HB2	2.17	0.44
2:B:86:ILE:O	2:B:86:ILE:HG22	2.17	0.44
2:A:110:ARG:HH12	2:A:112:VAL:HG22	1.80	0.44
2:B:48:LEU:HD12	2:B:188:LEU:HD13	1.99	0.44
2:B:139:LEU:HD13	2:B:220:LEU:CD2	2.47	0.44
2:B:22:HIS:CD2	2:B:23:MET:H	2.31	0.44
2:A:273:LEU:HD23	2:A:273:LEU:N	2.32	0.44
2:A:211:ARG:NE	2:A:219:LEU:HD23	2.33	0.44
2:A:57:THR:HG23	2:A:58:PRO:HD2	1.99	0.44
2:A:66:MET:SD	2:A:74:ILE:HD11	2.58	0.44
2:B:66:MET:HE2	2:B:98:LEU:O	2.18	0.44
2:A:136:LEU:HD12	2:A:136:LEU:H	1.79	0.44
2:B:165:GLN:N	2:B:166:PRO:CD	2.80	0.44
2:B:22:HIS:HB3	2:B:25:TRP:CD1	2.52	0.44
2:B:33:TRP:CG	2:B:256:ALA:HB2	2.52	0.44
2:A:237:GLN:HA	2:A:262:CYS:SG	2.57	0.44
2:A:269:LEU:O	2:A:271:TYR:N	2.51	0.44
2:A:49:ILE:HG22	2:A:50:ILE:N	2.32	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:258:LYS:NZ	2:B:258:LYS:CB	2.80	0.44
2:A:195:SER:N	2:A:196:PRO:CD	2.80	0.44
2:A:26:LEU:CD1	2:A:36:THR:OG1	2.66	0.44
2:B:61:ILE:HD11	2:B:117:GLU:CB	2.48	0.44
2:B:165:GLN:N	2:B:166:PRO:HD2	2.33	0.43
2:B:63:GLN:HE22	2:B:98:LEU:H	1.65	0.43
2:A:70:ALA:CB	2:A:109:TYR:CE2	3.01	0.43
2:A:120:LYS:NZ	2:A:121:ARG:HH12	2.17	0.43
2:A:123:GLU:CD	2:A:124:ASP:H	2.21	0.43
2:B:147:THR:HG23	2:B:261:ILE:HA	1.99	0.43
2:B:66:MET:HE2	2:B:100:PRO:CD	2.48	0.43
2:A:162:PRO:C	2:A:164:SER:N	2.70	0.43
2:B:66:MET:HE1	2:B:100:PRO:HD3	1.99	0.43
2:B:124:ASP:HA	2:B:193:LYS:HE3	2.00	0.43
2:A:121:ARG:NH1	2:A:121:ARG:HG3	2.32	0.43
2:A:225:ASP:HB3	2:A:268:TYR:HE1	1.83	0.43
2:B:43:VAL:HG21	2:B:44:ARG:NH1	2.26	0.43
2:B:199:PHE:HE2	2:B:224:ASP:HA	1.82	0.43
2:B:26:LEU:HA	2:B:26:LEU:HD13	1.81	0.43
2:B:79:GLN:O	2:B:80:ARG:C	2.56	0.43
2:B:97:PRO:HG2	2:B:114:ASP:HB3	1.99	0.43
1:E:9:DA:C4'	1:E:10:DT:OP1	2.55	0.43
2:A:94:TRP:CE3	2:A:118:VAL:HG22	2.53	0.43
2:A:263:GLN:NE2	2:A:264:LYS:H	2.17	0.43
2:A:26:LEU:HD11	2:A:36:THR:CG2	2.49	0.43
2:A:62:LYS:HA	2:A:62:LYS:HD3	1.86	0.43
2:A:89:PRO:HG2	2:B:162:PRO:HB3	2.00	0.43
2:B:72:LEU:O	2:B:73:GLY:C	2.57	0.43
2:A:203:LEU:HD23	2:A:223:VAL:O	2.19	0.42
2:B:26:LEU:HD11	2:B:36:THR:CG2	2.32	0.42
2:B:77:HIS:O	2:B:81:LEU:HD22	2.18	0.42
2:A:106:THR:OG1	2:A:108:ASP:HB2	2.19	0.42
2:A:67:SER:O	2:A:68:GLN:C	2.58	0.42
2:B:103:LYS:HA	2:B:103:LYS:HE3	2.00	0.42
2:A:230:ALA:HB1	2:A:235:ASP:HB2	2.00	0.42
2:A:269:LEU:C	2:A:271:TYR:N	2.71	0.42
2:B:248:GLY:O	2:B:250:LEU:N	2.52	0.42
2:A:139:LEU:HD23	2:A:139:LEU:O	2.20	0.42
2:A:68:GLN:NE2	2:A:69:GLU:N	2.68	0.42
2:A:70:ALA:CB	2:A:109:TYR:HE2	2.32	0.42
2:B:269:LEU:C	2:B:271:TYR:H	2.22	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:133:TYR:CD2	2:A:134:ASN:N	2.88	0.42
2:A:176:GLU:HG2	2:A:176:GLU:O	2.19	0.42
2:A:82:LEU:CD1	2:B:162:PRO:HG2	2.50	0.42
2:B:143:HIS:HA	2:B:231:THR:HA	2.01	0.42
2:A:165:GLN:H	2:A:165:GLN:HG3	1.46	0.41
2:A:30:PRO:O	2:A:37:GLY:HA3	2.20	0.41
2:B:71:ARG:HG2	2:B:71:ARG:HH11	1.85	0.41
2:B:145:TRP:CZ2	2:B:233:GLU:HB2	2.54	0.41
2:B:52:LEU:HD21	2:B:167:LEU:HD21	2.02	0.41
2:B:66:MET:CE	2:B:100:PRO:CD	2.98	0.41
2:B:69:GLU:H	2:B:69:GLU:CD	2.22	0.41
2:A:98:LEU:CD1	2:A:111:PRO:HB2	2.51	0.41
2:A:80:ARG:CG	2:A:81:LEU:HD22	2.48	0.41
2:B:143:HIS:O	2:B:146:TYR:HE1	2.04	0.41
2:B:248:GLY:C	2:B:250:LEU:H	2.23	0.41
2:B:72:LEU:O	2:B:75:LYS:HB3	2.20	0.41
1:E:3:DT:H3	2:A:67:SER:HG	1.66	0.41
2:B:125:ILE:O	2:B:126:HIS:C	2.58	0.41
2:B:151:LEU:CD2	2:B:254:ALA:HB2	2.50	0.41
1:E:5:DA:H1'	1:E:6:DT:H5''	2.03	0.41
2:A:103:LYS:O	2:A:106:THR:HG22	2.21	0.41
2:A:146:TYR:CD2	2:A:273:LEU:HD11	2.55	0.41
2:B:131:ASN:OD1	2:B:132:PRO:HD2	2.21	0.41
2:B:265:GLN:HA	2:B:273:LEU:O	2.21	0.41
2:A:268:TYR:CE1	2:A:269:LEU:HD13	2.55	0.41
2:B:94:TRP:CE2	2:B:167:LEU:HD12	2.56	0.41
1:E:5:DA:H1'	1:E:6:DT:C5'	2.51	0.41
2:A:158:LEU:HD12	2:A:158:LEU:N	2.35	0.41
2:B:80:ARG:NH1	2:B:80:ARG:HG2	2.36	0.41
2:A:268:TYR:O	2:A:271:TYR:HB2	2.21	0.41
2:B:32:ALA:HB2	2:B:244:LEU:O	2.22	0.40
2:B:81:LEU:HD12	2:B:86:ILE:HB	2.03	0.40
2:A:139:LEU:HD23	2:A:139:LEU:C	2.42	0.40
2:A:60:SER:HB2	2:A:93:PRO:O	2.20	0.40
2:B:272:LEU:C	2:B:273:LEU:HD23	2.42	0.40
1:E:10:DT:H2''	1:E:11:DG:OP2	2.20	0.40
2:A:218:ILE:HG22	2:A:220:LEU:HD12	2.03	0.40
2:A:40:GLY:N	2:A:253:ARG:HG3	2.36	0.40
2:B:212:ILE:H	2:B:212:ILE:HG13	1.55	0.40
2:A:94:TRP:HE3	2:A:118:VAL:HG22	1.87	0.40
2:A:158:LEU:CD1	2:A:158:LEU:N	2.84	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:211:ARG:CD	2:A:219:LEU:HB3	2.52	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	A	257/259 (99%)	215 (84%)	32 (12%)	10 (4%)	3	19
2	B	255/259 (98%)	201 (79%)	42 (16%)	12 (5%)	2	15
All	All	512/518 (99%)	416 (81%)	74 (14%)	22 (4%)	3	17

All (22) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	A	132	PRO
2	A	133	TYR
2	A	163	THR
2	B	72	LEU
2	B	207	LEU
2	A	54	ALA
2	A	223	VAL
2	B	38	GLY
2	B	134	ASN
2	B	249	ASN
2	B	277	GLN
2	A	37	GLY
2	B	172	TRP
2	B	223	VAL
2	A	176	GLU
2	A	107	ASN
2	A	130	PRO

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Mol	Chain	Res	Type
2	B	115	LEU
2	B	132	PRO
2	B	166	PRO
2	A	215	PRO
2	B	58	PRO

### 5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	A	227/227 (100%)	207 (91%)	20 (9%)	11	39
2	B	226/227 (100%)	200 (88%)	26 (12%)	6	25
All	All	453/454 (100%)	407 (90%)	46 (10%)	8	30

All (46) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	A	23	MET
2	A	26	LEU
2	A	31	GLN
2	A	44	ARG
2	A	65	PRO
2	A	82	LEU
2	A	98	LEU
2	A	111	PRO
2	A	116	ARG
2	A	127	PRO
2	A	160	LEU
2	A	165	GLN
2	A	173	ARG
2	A	204	HIS
2	A	211	ARG
2	A	257	LYS
2	A	261	ILE
2	A	267	LYS

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Mol	Chain	Res	Type
2	A	269	LEU
2	A	277	GLN
2	B	26	LEU
2	B	41	LEU
2	B	43	VAL
2	B	52	LEU
2	B	72	LEU
2	B	74	ILE
2	B	80	ARG
2	B	81	LEU
2	B	88	VAL
2	B	102	LYS
2	B	103	LYS
2	B	110	ARG
2	B	128	THR
2	B	149	LEU
2	B	163	THR
2	B	167	LEU
2	B	172	TRP
2	B	173	ARG
2	B	176	GLU
2	B	182	GLN
2	B	184	THR
2	B	212	ILE
2	B	225	ASP
2	B	235	ASP
2	B	274	LYS
2	B	278	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
2	A	31	GLN
2	A	68	GLN
2	A	113	GLN
2	A	161	HIS
2	A	165	GLN
2	A	194	ASN
2	A	237	GLN
2	A	245	GLN
2	A	249	ASN
2	A	263	GLN

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Mol	Chain	Res	Type
2	B	22	HIS
2	B	68	GLN
2	B	113	GLN
2	B	144	GLN
2	B	161	HIS
2	B	204	HIS
2	B	221	GLN
2	B	245	GLN
2	B	249	ASN
2	B	265	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](#)

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

## 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

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](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
2	A	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A	261:ILE	C	262:CYS	N	1.19

## 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	E	11/11 (100%)	0.37	1 (9%) 9 3	15, 40, 98, 98	0
1	F	11/11 (100%)	0.34	2 (18%) 1 0	15, 26, 99, 100	0
2	A	259/259 (100%)	-0.69	1 (0%) 92 78	3, 16, 44, 74	0
2	B	257/259 (99%)	-0.55	1 (0%) 92 78	5, 25, 64, 91	0
All	All	538/540 (99%)	-0.58	5 (0%) 84 62	3, 20, 55, 100	0

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	1	DT	4.5
2	B	106	THR	3.4
1	E	2	DT	3.0
2	A	104	PRO	2.4
1	F	2	DT	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 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.