



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 09:40 PM UTC

PDB ID : 3US0 / pdb_00003us0
Title : Structure of p63 DNA Binding Domain in Complex with a 22 Base Pair A/T Rich Response Element Containing a Two Base Pair "AT" Spacer Between Half Sites
Authors : Chen, C.; Herzberg, O.
Deposited on : 2011-11-22
Resolution : 2.50 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 ⓘ 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 ⓘ](#)) were used in the production of this report:

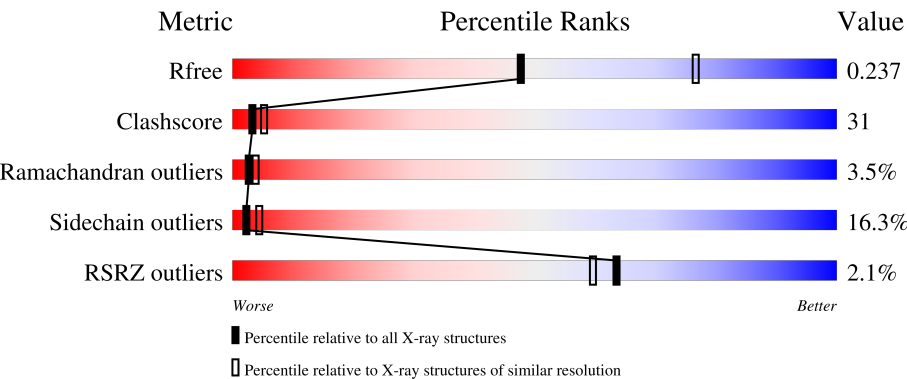
MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

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 2.50 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}	180053	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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 $\leq 5\%$. 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	203	2% 37% 42% 17% .
1	B	203	% 46% 41% 7% 5%
1	C	203	4% 42% 40% 13% .
1	D	203	% 53% 34% 7% 5%
2	E	22	27% 73%

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Mol	Chain	Length	Quality of chain
2	F	22	32% 64% 5%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7279 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 Tumor protein 63.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	195	Total	C	N	O	S	0	0	0
			1503	940	266	285	12			
1	B	193	Total	C	N	O	S	0	0	0
			1491	936	262	280	13			
1	C	196	Total	C	N	O	S	0	0	0
			1509	946	263	288	12			
1	D	192	Total	C	N	O	S	0	0	0
			1482	931	258	280	13			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	121	GLY	-	expression tag	UNP Q9H3D4
A	122	SER	-	expression tag	UNP Q9H3D4
A	123	HIS	-	expression tag	UNP Q9H3D4
A	124	MET	-	expression tag	UNP Q9H3D4
A	125	ALA	-	expression tag	UNP Q9H3D4
A	126	SER	-	expression tag	UNP Q9H3D4
B	121	GLY	-	expression tag	UNP Q9H3D4
B	122	SER	-	expression tag	UNP Q9H3D4
B	123	HIS	-	expression tag	UNP Q9H3D4
B	124	MET	-	expression tag	UNP Q9H3D4
B	125	ALA	-	expression tag	UNP Q9H3D4
B	126	SER	-	expression tag	UNP Q9H3D4
C	121	GLY	-	expression tag	UNP Q9H3D4
C	122	SER	-	expression tag	UNP Q9H3D4
C	123	HIS	-	expression tag	UNP Q9H3D4
C	124	MET	-	expression tag	UNP Q9H3D4
C	125	ALA	-	expression tag	UNP Q9H3D4
C	126	SER	-	expression tag	UNP Q9H3D4
D	121	GLY	-	expression tag	UNP Q9H3D4
D	122	SER	-	expression tag	UNP Q9H3D4
D	123	HIS	-	expression tag	UNP Q9H3D4

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Chain	Residue	Modelled	Actual	Comment	Reference
D	124	MET	-	expression tag	UNP Q9H3D4
D	125	ALA	-	expression tag	UNP Q9H3D4
D	126	SER	-	expression tag	UNP Q9H3D4

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	22	Total	C	N	O	P	0	0	0
			448	218	79	130	21			
2	F	22	Total	C	N	O	P	0	0	0
			448	218	79	130	21			

- Molecule 3 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Zn	0	0
			1	1		
3	B	1	Total	Zn	0	0
			1	1		
3	C	1	Total	Zn	0	0
			1	1		
3	D	1	Total	Zn	0	0
			1	1		

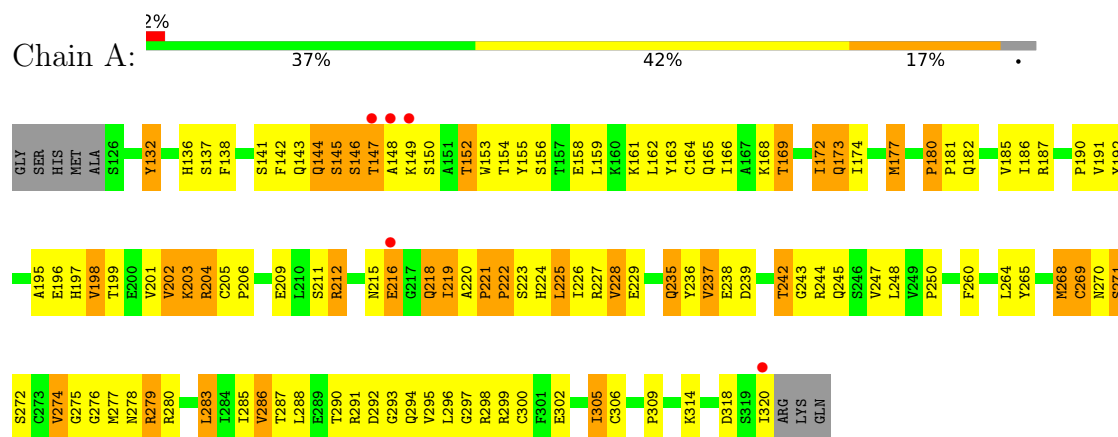
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	71	Total	O	0	0
			71	71		
4	B	105	Total	O	0	0
			105	105		
4	C	65	Total	O	0	0
			65	65		
4	D	103	Total	O	0	0
			103	103		
4	E	24	Total	O	0	0
			24	24		
4	F	26	Total	O	0	0
			26	26		

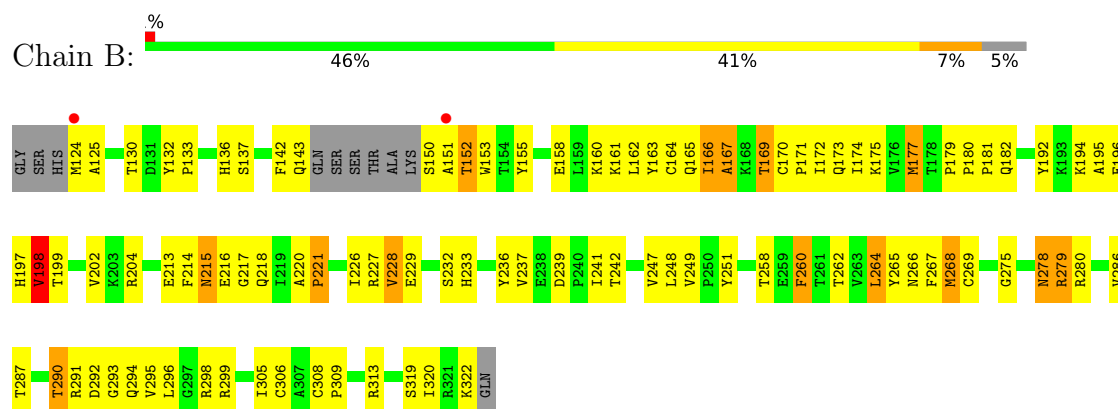
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

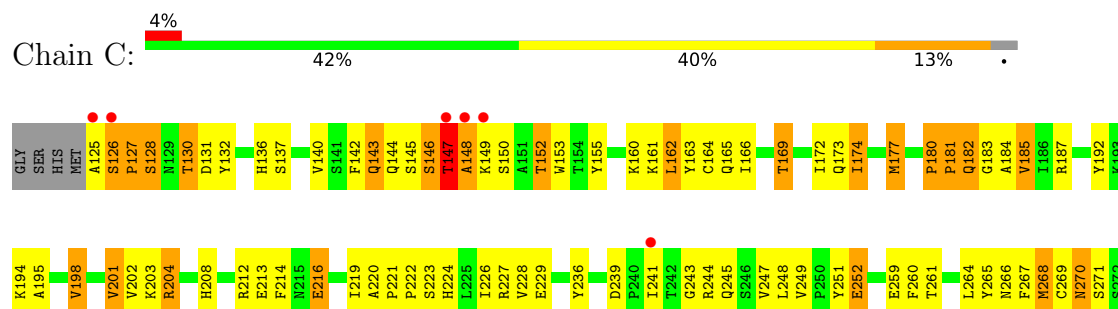
• Molecule 1: Tumor protein 63

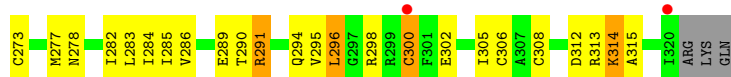


• Molecule 1: Tumor protein 63



• Molecule 1: Tumor protein 63





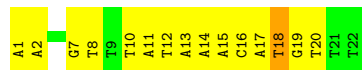
• Molecule 1: Tumor protein 63



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



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



4 Data and refinement statistics

Property	Value	Source
Space group	P 63	Depositor
Cell constants a, b, c, α , β , γ	141.81Å 141.81Å 119.60Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	59.80 – 2.50 59.80 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.9 (59.80-2.50) 100.0 (59.80-2.50)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.44 (at 2.40Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7_650)	Depositor
R, R_{free}	0.206 , 0.239 0.204 , 0.237	Depositor DCC
R_{free} test set	2728 reflections (5.11%)	wwPDB-VP
Wilson B-factor (Å ²)	47.0	Xtriage
Anisotropy	0.408	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 48.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.052 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	7279	wwPDB-VP
Average B, all atoms (Å ²)	67.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.82% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

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

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# $ Z > 5$	RMSZ	# $ Z > 5$
1	A	0.71	0/1539	1.18	12/2096 (0.6%)
1	B	0.70	0/1526	1.13	9/2074 (0.4%)
1	C	0.67	0/1545	1.19	16/2104 (0.8%)
1	D	0.69	0/1517	1.16	10/2064 (0.5%)
2	E	0.31	0/502	0.99	0/773
2	F	0.28	0/502	1.06	1/773 (0.1%)
All	All	0.65	0/7131	1.15	48/9884 (0.5%)

There are no bond length outliers.

All (48) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	268	MET	N-CA-C	8.94	123.79	113.15
1	C	126	SER	CA-C-N	8.55	130.53	119.84
1	C	126	SER	C-N-CA	8.55	130.53	119.84
1	B	268	MET	N-CA-C	8.41	122.31	112.72
1	D	268	MET	N-CA-C	8.14	122.79	113.02
1	C	180	PRO	CA-C-N	7.77	129.55	119.84
1	C	180	PRO	C-N-CA	7.77	129.55	119.84
1	C	268	MET	N-CA-C	7.66	122.27	113.15
1	A	293	GLY	N-CA-C	7.58	125.05	115.42
1	D	132	TYR	CA-C-N	7.58	127.12	119.24
1	D	132	TYR	C-N-CA	7.58	127.12	119.24
1	B	198	VAL	CB-CA-C	-7.54	102.16	112.04
2	F	18	DT	P-O5'-C5'	-7.21	109.19	120.00
1	D	226	ILE	CB-CA-C	-7.18	102.00	110.91
1	A	180	PRO	CA-C-N	7.06	128.66	119.84
1	A	180	PRO	C-N-CA	7.06	128.66	119.84
1	B	226	ILE	CB-CA-C	-6.79	102.50	110.91
1	D	275	GLY	N-CA-C	-6.64	106.79	114.69

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	132	TYR	CA-C-N	6.62	127.15	119.47
1	C	132	TYR	C-N-CA	6.62	127.15	119.47
1	A	205	CYS	CA-C-N	6.35	125.84	119.24
1	A	205	CYS	C-N-CA	6.35	125.84	119.24
1	A	226	ILE	N-CA-C	6.20	116.79	107.80
1	D	212	ARG	N-CA-C	-6.12	104.69	111.36
1	B	221	PRO	N-CA-C	-6.10	103.26	110.70
1	B	132	TYR	CA-C-N	6.05	125.77	119.05
1	B	132	TYR	C-N-CA	6.05	125.77	119.05
1	C	164	CYS	N-CA-C	5.95	118.20	108.55
1	A	221	PRO	CA-C-N	5.84	127.14	119.84
1	A	221	PRO	C-N-CA	5.84	127.14	119.84
1	A	237	VAL	N-CA-C	5.81	115.90	107.88
1	B	293	GLY	N-CA-C	5.79	122.83	114.10
1	C	261	THR	N-CA-C	-5.69	100.40	109.96
1	D	173	GLN	N-CA-C	5.65	117.50	108.41
1	A	132	TYR	CA-C-N	5.47	125.12	119.05
1	A	132	TYR	C-N-CA	5.47	125.12	119.05
1	D	174	ILE	CB-CA-C	-5.43	103.77	111.21
1	D	198	VAL	CB-CA-C	-5.41	104.95	112.04
1	B	275	GLY	N-CA-C	-5.30	108.38	114.69
1	C	128	SER	N-CA-C	5.28	117.36	109.59
1	D	178	THR	N-CA-C	-5.23	102.19	108.25
1	B	260	PHE	N-CA-C	5.18	117.51	108.76
1	C	212	ARG	N-CA-C	-5.14	105.37	111.69
1	C	270	ASN	N-CA-C	-5.12	104.11	110.41
1	C	185	VAL	N-CA-C	5.11	116.26	108.80
1	C	174	ILE	CB-CA-C	-5.10	103.06	110.81
1	C	252	GLU	CA-C-N	-5.04	115.19	120.38
1	C	252	GLU	C-N-CA	-5.04	115.19	120.38

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	1503	0	1455	111	0
1	B	1491	0	1454	77	0
1	C	1509	0	1468	96	0
1	D	1482	0	1434	73	0
2	E	448	0	253	28	0
2	F	448	0	253	18	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	A	71	0	0	4	0
4	B	105	0	0	14	0
4	C	65	0	0	6	0
4	D	103	0	0	5	0
4	E	24	0	0	5	0
4	F	26	0	0	5	0
All	All	7279	0	6317	403	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 31.

All (403) 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:203:LYS:HB3	1:A:245:GLN:HE22	1.11	1.08
1:C:182:GLN:HE21	1:C:182:GLN:HA	1.20	1.04
2:F:12:DT:H2''	2:F:13:DA:H5'	1.44	0.99
2:F:11:DA:OP2	4:F:373:HOH:O	1.81	0.98
1:C:285:ILE:HG12	1:C:300:CYS:SG	2.08	0.93
2:E:11:DA:N7	4:E:369:HOH:O	2.02	0.91
1:B:290:THR:HG22	1:B:294:GLN:H	1.36	0.90
1:D:290:THR:HG22	1:D:294:GLN:H	1.37	0.90
1:A:204:ARG:HD2	1:A:268:MET:HB2	1.53	0.89
1:A:137:SER:HB2	1:A:177:MET:HB2	1.54	0.88
1:B:214:PHE:C	1:B:215:ASN:HD22	1.82	0.87
1:B:133:PRO:O	4:B:368:HOH:O	1.91	0.86
1:D:176:VAL:O	1:D:177:MET:HE3	1.75	0.86
1:D:279:ARG:HG3	1:D:279:ARG:HH11	1.39	0.86
1:B:169:THR:HB	1:B:229:GLU:OE2	1.76	0.85
1:B:215:ASN:HD22	1:B:215:ASN:N	1.72	0.85
1:D:227:ARG:NH1	1:D:268:MET:HE2	1.92	0.85
1:B:199:THR:OG1	4:B:392:HOH:O	1.95	0.84

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:203:LYS:HB3	1:A:245:GLN:NE2	1.92	0.84
1:B:158:GLU:HG3	4:B:345:HOH:O	1.76	0.83
1:C:204:ARG:HD2	1:C:268:MET:HB2	1.60	0.83
1:D:278:ASN:ND2	1:D:279:ARG:HH12	1.76	0.82
1:B:177:MET:HE3	1:B:177:MET:HA	1.62	0.82
1:C:187:ARG:HB2	1:C:248:LEU:CD2	2.09	0.82
1:A:142:PHE:CE1	1:A:172:ILE:HD13	2.16	0.81
1:D:290:THR:HG23	1:D:292:ASP:H	1.45	0.80
1:C:125:ALA:HA	4:C:364:HOH:O	1.82	0.80
1:D:204:ARG:HD2	1:D:268:MET:HB2	1.62	0.80
2:F:10:DT:O2	4:F:371:HOH:O	1.99	0.79
1:A:212:ARG:CG	1:A:212:ARG:HH21	1.94	0.79
1:A:177:MET:HE3	1:A:177:MET:HA	1.65	0.79
1:A:192:TYR:CE2	1:A:202:VAL:HG13	2.19	0.78
1:A:204:ARG:HD2	1:A:268:MET:CB	2.13	0.78
1:D:278:ASN:HD22	1:D:279:ARG:HH12	1.31	0.78
1:C:177:MET:HE3	1:C:177:MET:HA	1.67	0.77
1:D:204:ARG:NH1	1:D:268:MET:HB3	1.99	0.77
1:D:177:MET:HE3	1:D:177:MET:HA	1.67	0.77
1:D:278:ASN:HD22	1:D:279:ARG:NH1	1.83	0.77
1:A:187:ARG:HB2	1:A:248:LEU:CD2	2.15	0.77
1:A:136:HIS:CE1	1:A:180:PRO:HA	2.20	0.77
2:F:12:DT:H73	4:F:36:HOH:O	1.86	0.76
1:C:169:THR:HB	1:C:229:GLU:OE2	1.86	0.75
1:A:173:GLN:HG2	1:A:260:PHE:CG	2.21	0.75
2:E:12:DT:H73	4:E:104:HOH:O	1.85	0.75
1:A:166:ILE:O	4:A:383:HOH:O	2.04	0.75
1:B:143:GLN:HB3	1:B:173:GLN:HE22	1.52	0.75
1:B:169:THR:HA	1:B:266:ASN:ND2	2.03	0.74
1:C:187:ARG:HB2	1:C:248:LEU:HD22	1.70	0.73
1:B:290:THR:HG23	1:B:292:ASP:H	1.52	0.73
1:D:227:ARG:HH11	1:D:268:MET:HE2	1.54	0.73
1:C:183:GLY:HA3	1:C:291:ARG:HB2	1.70	0.73
1:B:204:ARG:NH1	1:B:268:MET:HB3	2.04	0.73
2:F:12:DT:C2'	2:F:13:DA:H5'	2.18	0.72
1:A:290:THR:HG22	1:A:294:GLN:H	1.52	0.72
1:B:143:GLN:OE1	4:B:380:HOH:O	2.06	0.72
1:B:320:ILE:C	1:B:322:LYS:H	1.96	0.72
1:B:152:THR:HG22	1:B:153:TRP:CD1	2.24	0.72
1:D:142:PHE:CE1	1:D:172:ILE:HG12	2.24	0.72
1:D:187:ARG:HB3	1:D:287:THR:CG2	2.20	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:143:GLN:HB3	1:B:173:GLN:NE2	2.05	0.72
1:B:204:ARG:HD2	1:B:268:MET:HB2	1.72	0.72
1:D:182:GLN:HA	1:D:182:GLN:NE2	2.05	0.71
2:E:19:DG:H2''	2:E:20:DT:C5'	2.20	0.71
1:A:219:ILE:HG22	1:A:219:ILE:O	1.90	0.71
2:E:17:DA:H2''	2:E:18:DT:H5'	1.73	0.70
1:A:187:ARG:HB2	1:A:248:LEU:HD22	1.72	0.70
1:B:152:THR:HG22	1:B:153:TRP:HD1	1.56	0.70
1:A:218:GLN:HE21	1:A:218:GLN:HA	1.56	0.69
1:D:142:PHE:CE2	1:D:155:TYR:CE1	2.81	0.69
1:D:279:ARG:HG3	1:D:279:ARG:NH1	2.04	0.69
1:B:298:ARG:O	1:B:299:ARG:HD2	1.92	0.69
1:D:169:THR:HA	1:D:266:ASN:ND2	2.09	0.68
1:C:290:THR:HG22	1:C:294:GLN:H	1.58	0.68
1:C:144:GLN:OE1	1:C:144:GLN:HA	1.92	0.68
2:E:19:DG:H2''	2:E:20:DT:H5''	1.75	0.68
2:F:11:DA:H3'	4:F:374:HOH:O	1.94	0.67
1:C:182:GLN:HB3	1:C:291:ARG:NH2	2.10	0.67
1:C:182:GLN:HA	1:C:182:GLN:NE2	2.02	0.67
1:C:187:ARG:CZ	1:C:248:LEU:HD21	2.25	0.67
1:C:163:TYR:CE2	1:C:313:ARG:HA	2.30	0.66
1:C:290:THR:CG2	1:C:294:GLN:H	2.08	0.66
1:C:181:PRO:HG2	1:C:296:LEU:HD21	1.76	0.66
1:C:290:THR:CG2	1:C:294:GLN:HB3	2.25	0.66
1:A:212:ARG:HH21	1:A:212:ARG:HG2	1.60	0.66
1:A:172:ILE:HG13	1:A:265:TYR:HD2	1.59	0.66
1:A:271:SER:O	1:A:279:ARG:HA	1.96	0.66
1:A:290:THR:HG22	1:A:294:GLN:N	2.10	0.65
1:C:204:ARG:NH1	1:C:208:HIS:HB3	2.12	0.65
1:B:233:HIS:ND1	4:B:325:HOH:O	2.29	0.65
1:C:187:ARG:HB2	1:C:248:LEU:HD23	1.76	0.65
2:E:8:DT:H2''	2:E:9:DT:H5''	1.77	0.65
1:A:239:ASP:O	1:A:243:GLY:N	2.24	0.65
1:D:187:ARG:HB3	1:D:287:THR:HG23	1.78	0.65
1:B:137:SER:HB2	1:B:177:MET:HB2	1.79	0.65
1:D:142:PHE:CE2	1:D:155:TYR:CD1	2.85	0.65
1:D:177:MET:HA	1:D:177:MET:CE	2.27	0.65
1:C:146:SER:O	1:C:148:ALA:N	2.31	0.64
1:D:316:ASP:O	1:D:320:ILE:HG23	1.96	0.64
2:E:17:DA:N3	4:E:113:HOH:O	2.30	0.64
1:C:213:GLU:O	1:C:214:PHE:HB2	1.98	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:169:THR:HB	1:A:229:GLU:OE2	1.98	0.64
1:C:161:LYS:HG3	1:C:302:GLU:HB3	1.80	0.64
1:C:166:ILE:HB	1:C:270:ASN:OD1	1.97	0.64
1:A:166:ILE:HA	1:A:305:ILE:HD12	1.78	0.64
1:B:278:ASN:O	1:B:279:ARG:HG2	1.97	0.64
1:A:195:ALA:O	1:A:198:VAL:HG22	1.97	0.63
1:B:215:ASN:N	1:B:215:ASN:ND2	2.46	0.63
1:D:290:THR:HG22	1:D:294:GLN:N	2.10	0.63
1:A:187:ARG:CZ	1:A:248:LEU:HD21	2.29	0.63
1:C:172:ILE:HD12	1:C:265:TYR:CD2	2.34	0.63
1:C:226:ILE:HD11	1:C:284:ILE:HD12	1.80	0.63
1:D:320:ILE:C	1:D:322:LYS:H	2.06	0.63
1:D:204:ARG:HH11	1:D:268:MET:HB3	1.64	0.62
1:C:290:THR:HG22	1:C:294:GLN:N	2.14	0.62
1:C:290:THR:HG21	1:C:294:GLN:HB3	1.82	0.62
1:C:228:VAL:HG12	1:C:247:VAL:HG21	1.81	0.62
1:A:152:THR:OG1	1:A:168:LYS:HG3	2.00	0.62
1:D:224:HIS:CE1	1:D:245:GLN:HB3	2.34	0.62
1:C:147:THR:O	1:C:149:LYS:N	2.33	0.62
1:A:215:ASN:O	1:A:216:GLU:C	2.43	0.61
1:C:277:MET:HE3	1:C:282:ILE:HG23	1.83	0.61
1:D:136:HIS:CE1	1:D:180:PRO:HA	2.35	0.61
1:C:271:SER:HB2	4:C:333:HOH:O	2.01	0.61
1:B:204:ARG:HH11	1:B:268:MET:HB3	1.65	0.61
1:A:142:PHE:HE1	1:A:172:ILE:HD13	1.62	0.60
1:A:144:GLN:OE1	1:A:144:GLN:HA	2.02	0.60
1:C:182:GLN:HB3	1:C:291:ARG:HH21	1.66	0.60
1:B:152:THR:CG2	1:B:153:TRP:HD1	2.15	0.60
1:A:235:GLN:HG3	1:A:235:GLN:O	2.01	0.60
1:B:228:VAL:HG12	1:B:247:VAL:HG21	1.84	0.59
2:F:17:DA:H2''	2:F:18:DT:H5'	1.84	0.59
1:B:169:THR:HG22	4:B:342:HOH:O	2.01	0.59
1:D:290:THR:HG23	1:D:292:ASP:N	2.14	0.59
1:C:194:LYS:O	1:C:198:VAL:HG13	2.03	0.59
1:A:201:VAL:HG12	1:A:202:VAL:N	2.18	0.58
1:A:161:LYS:HD3	1:A:163:TYR:CZ	2.38	0.58
1:A:142:PHE:CE2	1:A:155:TYR:CE1	2.91	0.58
1:C:192:TYR:CE2	1:C:202:VAL:HG22	2.38	0.58
2:E:8:DT:C2'	2:E:9:DT:H5''	2.34	0.58
1:C:182:GLN:HE21	1:C:182:GLN:CA	1.99	0.58
1:B:163:TYR:CE2	1:B:313:ARG:HA	2.39	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:290:THR:CG2	1:A:294:GLN:H	2.17	0.58
1:B:142:PHE:CE2	1:B:155:TYR:CE1	2.91	0.58
2:E:15:DA:H1'	2:E:16:DC:H5''	1.85	0.58
1:A:285:ILE:HG12	1:A:300:CYS:SG	2.44	0.57
1:C:224:HIS:CE1	1:C:236:TYR:HB3	2.39	0.57
1:D:153:TRP:NE1	1:D:171:PRO:HD2	2.20	0.57
1:D:130:THR:HG23	4:D:335:HOH:O	2.05	0.56
1:C:203:LYS:O	4:C:5:HOH:O	2.17	0.56
1:D:201:VAL:HG12	1:D:244:ARG:HG2	1.86	0.56
1:D:228:VAL:HG12	1:D:247:VAL:HG21	1.88	0.56
1:D:318:ASP:O	1:D:322:LYS:HG3	2.06	0.56
2:E:8:DT:H2''	2:E:9:DT:C5'	2.35	0.56
1:B:150:SER:HA	4:B:116:HOH:O	2.04	0.56
1:B:217:GLY:O	1:B:218:GLN:HG2	2.06	0.56
1:D:298:ARG:O	1:D:299:ARG:HD2	2.05	0.56
1:B:204:ARG:NH1	1:B:268:MET:O	2.39	0.55
1:C:169:THR:HA	1:C:266:ASN:ND2	2.21	0.55
1:C:162:LEU:HD22	1:C:163:TYR:N	2.21	0.55
1:D:169:THR:HA	1:D:266:ASN:HD21	1.71	0.55
1:A:185:VAL:HG12	1:A:250:PRO:HA	1.87	0.55
1:A:221:PRO:HG2	1:A:224:HIS:CG	2.42	0.55
1:A:228:VAL:HG12	1:A:247:VAL:HG21	1.89	0.55
1:A:269:CYS:O	1:A:305:ILE:HG12	2.06	0.55
1:B:169:THR:HA	1:B:266:ASN:HD21	1.69	0.55
2:E:1:DA:HO5'	2:E:1:DA:H8	1.53	0.55
2:F:19:DG:H2''	2:F:20:DT:H5'	1.89	0.55
1:D:290:THR:CG2	1:D:294:GLN:H	2.13	0.54
1:D:171:PRO:HA	1:D:264:LEU:HD12	1.89	0.54
1:C:267:PHE:HB3	1:C:305:ILE:CD1	2.37	0.54
1:B:150:SER:HB2	4:B:55:HOH:O	2.07	0.54
1:D:279:ARG:HH11	1:D:279:ARG:CG	2.13	0.54
2:F:12:DT:OP2	4:F:374:HOH:O	2.18	0.54
1:A:172:ILE:HG13	1:A:265:TYR:CD2	2.42	0.54
1:C:172:ILE:HD12	1:C:265:TYR:HD2	1.70	0.54
1:A:202:VAL:HG23	4:A:342:HOH:O	2.07	0.54
1:D:204:ARG:NH1	1:D:268:MET:O	2.40	0.54
1:B:177:MET:HA	1:B:177:MET:CE	2.36	0.53
1:B:143:GLN:N	4:B:380:HOH:O	1.87	0.53
1:D:169:THR:HB	1:D:229:GLU:OE2	2.08	0.53
1:B:320:ILE:C	1:B:322:LYS:N	2.66	0.53
1:C:204:ARG:HB2	1:C:223:SER:O	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:136:HIS:CE1	1:B:180:PRO:HA	2.43	0.53
1:A:146:SER:C	1:A:148:ALA:H	2.16	0.53
1:A:201:VAL:CG1	1:A:202:VAL:N	2.71	0.53
1:A:212:ARG:HH21	1:A:212:ARG:HG3	1.70	0.53
1:D:320:ILE:C	1:D:322:LYS:N	2.66	0.53
1:A:196:GLU:CD	1:A:196:GLU:H	2.16	0.52
1:B:279:ARG:HH21	1:B:279:ARG:CG	2.22	0.52
2:E:16:DC:H6	2:E:16:DC:H5'	1.74	0.52
2:F:13:DA:H1'	2:F:14:DA:H5''	1.90	0.52
2:E:8:DT:H1'	2:E:9:DT:H5''	1.91	0.52
2:E:11:DA:H1'	2:E:12:DT:H5'	1.92	0.52
1:D:187:ARG:HG3	1:D:248:LEU:HD23	1.91	0.52
1:D:204:ARG:HG2	1:D:269:CYS:SG	2.50	0.52
1:A:145:SER:HB3	1:A:154:THR:HA	1.91	0.52
1:C:147:THR:C	1:C:149:LYS:H	2.18	0.52
1:A:166:ILE:HA	1:A:305:ILE:CD1	2.40	0.52
1:B:143:GLN:CG	4:B:380:HOH:O	2.57	0.52
1:B:279:ARG:HH21	1:B:279:ARG:HG3	1.75	0.52
1:A:295:VAL:HG21	1:A:298:ARG:HH21	1.75	0.52
1:D:152:THR:N	4:D:392:HOH:O	2.30	0.51
2:E:11:DA:N6	4:E:369:HOH:O	2.17	0.51
1:B:279:ARG:CG	1:B:279:ARG:NH2	2.74	0.51
1:D:124:MET:N	4:D:19:HOH:O	2.42	0.51
1:C:145:SER:O	1:C:147:THR:N	2.43	0.51
1:C:201:VAL:HG12	1:C:244:ARG:HG2	1.91	0.51
1:C:267:PHE:HB3	1:C:305:ILE:HD13	1.92	0.51
1:A:138:PHE:O	1:A:299:ARG:NH1	2.43	0.51
1:A:166:ILE:HG22	4:A:332:HOH:O	2.11	0.51
1:D:142:PHE:CD2	1:D:155:TYR:CD1	2.98	0.51
1:A:187:ARG:CB	1:A:248:LEU:CD2	2.87	0.50
1:A:136:HIS:O	1:A:137:SER:C	2.54	0.50
1:C:184:ALA:HA	1:C:289:GLU:O	2.11	0.50
1:B:287:THR:HG22	4:B:93:HOH:O	2.12	0.50
1:A:218:GLN:HE21	1:A:218:GLN:CA	2.25	0.50
1:C:165:GLN:HA	1:C:306:CYS:O	2.11	0.50
1:B:171:PRO:HA	1:B:264:LEU:HD12	1.94	0.50
1:B:213:GLU:O	1:B:214:PHE:HB2	2.12	0.50
2:E:17:DA:H2''	2:E:18:DT:C5'	2.42	0.50
2:F:20:DT:H5'	2:F:20:DT:H6	1.77	0.50
1:A:165:GLN:HA	1:A:306:CYS:O	2.12	0.49
1:B:172:ILE:HD12	1:B:265:TYR:CD2	2.47	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:252:GLU:N	4:C:12:HOH:O	2.43	0.49
1:A:138:PHE:CE1	1:A:288:LEU:HB2	2.47	0.49
1:A:305:ILE:HG13	1:A:305:ILE:O	2.11	0.49
1:C:290:THR:HG22	1:C:294:GLN:HB3	1.93	0.49
1:A:161:LYS:HE3	1:A:302:GLU:OE2	2.12	0.49
1:B:214:PHE:C	1:B:215:ASN:ND2	2.63	0.49
1:A:295:VAL:O	1:A:295:VAL:HG13	2.12	0.49
1:C:148:ALA:O	1:C:149:LYS:HB2	2.12	0.49
1:A:173:GLN:HG2	1:A:260:PHE:CB	2.42	0.49
1:B:124:MET:N	4:B:334:HOH:O	2.44	0.49
1:C:177:MET:HA	1:C:177:MET:CE	2.40	0.49
1:C:195:ALA:HA	1:C:198:VAL:HG13	1.93	0.49
2:F:19:DG:H2''	2:F:20:DT:C5'	2.43	0.49
1:D:204:ARG:HH11	1:D:268:MET:CB	2.25	0.49
1:C:142:PHE:CE2	1:C:155:TYR:CD1	3.00	0.48
1:A:164:CYS:O	1:A:305:ILE:HA	2.13	0.48
1:A:148:ALA:O	1:A:150:SER:N	2.46	0.48
1:A:142:PHE:CE2	1:A:155:TYR:CD1	3.01	0.48
1:A:177:MET:HA	1:A:177:MET:CE	2.41	0.48
1:B:142:PHE:CE1	1:B:172:ILE:HG12	2.48	0.48
1:D:182:GLN:C	1:D:184:ALA:H	2.22	0.48
1:A:225:LEU:HD13	1:A:277:MET:HE2	1.94	0.48
1:A:290:THR:HG23	1:A:292:ASP:H	1.78	0.48
2:E:12:DT:H1'	2:E:13:DA:H5'	1.96	0.48
1:A:212:ARG:CG	1:A:212:ARG:NH2	2.63	0.48
1:C:137:SER:HB2	1:C:177:MET:HB2	1.95	0.48
1:D:174:ILE:HG13	1:D:186:ILE:CD1	2.44	0.47
1:A:152:THR:CG2	1:A:153:TRP:HD1	2.27	0.47
1:A:161:LYS:HD3	1:A:163:TYR:OH	2.15	0.47
1:D:166:ILE:C	1:D:168:LYS:H	2.22	0.47
1:D:182:GLN:HA	1:D:182:GLN:HE21	1.79	0.47
2:F:1:DA:H2''	2:F:2:DA:C8	2.49	0.47
1:C:251:TYR:CD2	1:C:251:TYR:C	2.93	0.47
1:A:209:GLU:HG3	1:A:222:PRO:HB2	1.97	0.47
1:C:283:LEU:HD23	1:C:302:GLU:HG3	1.97	0.47
1:D:221:PRO:HB2	1:D:224:HIS:CD2	2.49	0.47
1:A:141:SER:O	1:A:173:GLN:HB2	2.14	0.47
1:A:197:HIS:O	1:A:199:THR:N	2.48	0.47
1:D:174:ILE:HG13	1:D:186:ILE:HD13	1.96	0.47
1:A:143:GLN:HB3	1:A:173:GLN:OE1	2.15	0.47
1:B:278:ASN:C	1:B:279:ARG:HG2	2.40	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:136:HIS:CE1	1:C:180:PRO:HA	2.51	0.46
2:E:11:DA:C5	4:E:369:HOH:O	2.58	0.46
1:A:218:GLN:C	1:A:220:ALA:H	2.23	0.46
1:C:204:ARG:HG3	1:C:269:CYS:SG	2.55	0.46
1:C:192:TYR:CD2	1:C:202:VAL:HG22	2.50	0.46
1:D:228:VAL:HG23	1:D:264:LEU:O	2.16	0.46
1:B:179:PRO:HA	1:B:180:PRO:HD2	1.83	0.46
1:C:148:ALA:C	1:C:150:SER:H	2.24	0.46
1:C:183:GLY:O	1:C:185:VAL:HG13	2.16	0.46
1:A:204:ARG:HD2	1:A:268:MET:HB3	1.95	0.46
1:B:239:ASP:HB3	1:B:242:THR:OG1	2.16	0.46
2:E:22:DT:H5'	2:E:22:DT:C6	2.50	0.46
1:A:224:HIS:CE1	1:A:236:TYR:HB3	2.51	0.46
2:E:17:DA:H1'	2:E:18:DT:H5''	1.98	0.45
1:C:126:SER:HA	1:C:127:PRO:HD2	1.62	0.45
1:C:142:PHE:CE2	1:C:155:TYR:CE1	3.05	0.45
2:E:19:DG:H2''	2:E:20:DT:H5'	1.96	0.45
1:B:192:TYR:CE2	1:B:202:VAL:HG22	2.51	0.45
1:C:271:SER:HA	1:C:277:MET:HE2	1.97	0.45
1:D:163:TYR:CE2	1:D:313:ARG:HA	2.52	0.45
2:E:15:DA:C2'	2:E:16:DC:H5''	2.47	0.45
1:D:142:PHE:HE2	1:D:155:TYR:CE1	2.31	0.45
1:A:152:THR:HG22	1:A:153:TRP:HD1	1.80	0.45
1:B:174:ILE:HG22	1:B:175:LYS:N	2.31	0.45
1:C:182:GLN:NE2	1:C:182:GLN:CA	2.71	0.45
1:A:204:ARG:HG3	1:A:269:CYS:SG	2.56	0.45
1:D:227:ARG:HD2	1:D:268:MET:CG	2.47	0.45
1:A:173:GLN:HG2	1:A:260:PHE:CD2	2.52	0.45
1:A:270:ASN:C	1:A:272:SER:H	2.25	0.45
1:A:305:ILE:O	1:A:305:ILE:CG1	2.64	0.45
1:C:152:THR:CG2	1:C:153:TRP:HD1	2.29	0.45
1:C:204:ARG:NH1	1:C:268:MET:O	2.49	0.45
1:A:212:ARG:HG3	1:A:212:ARG:NH2	2.32	0.45
1:C:140:VAL:HG12	1:C:174:ILE:CD1	2.47	0.44
1:D:169:THR:CB	1:D:229:GLU:OE2	2.65	0.44
2:E:20:DT:H2'	2:E:21:DT:H72	1.99	0.44
1:A:145:SER:O	1:A:146:SER:C	2.60	0.44
1:A:238:GLU:HA	1:A:244:ARG:O	2.17	0.44
1:C:152:THR:HG22	1:C:153:TRP:HD1	1.81	0.44
1:C:187:ARG:CB	1:C:248:LEU:HD23	2.47	0.44
1:D:221:PRO:HB2	1:D:224:HIS:HD2	1.82	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:227:ARG:CZ	1:A:268:MET:HE2	2.48	0.44
1:A:227:ARG:NH1	1:A:268:MET:HE2	2.32	0.44
2:F:7:DG:H1'	2:F:8:DT:H5''	1.99	0.44
1:B:258:THR:HG22	1:B:260:PHE:H	1.82	0.44
1:D:142:PHE:HA	4:D:107:HOH:O	2.17	0.44
1:A:277:MET:O	1:A:280:ARG:HB2	2.18	0.44
1:B:290:THR:HG23	1:B:292:ASP:N	2.27	0.44
1:D:204:ARG:HB2	1:D:223:SER:O	2.17	0.44
1:A:142:PHE:CD2	1:A:155:TYR:CG	3.05	0.44
1:B:228:VAL:HA	1:B:264:LEU:O	2.18	0.44
2:F:15:DA:H1'	2:F:16:DC:H5''	1.99	0.44
1:B:220:ALA:O	1:B:221:PRO:C	2.61	0.44
1:D:164:CYS:O	1:D:305:ILE:HA	2.17	0.44
1:C:220:ALA:HB2	1:C:236:TYR:CZ	2.52	0.44
1:B:267:PHE:HB3	1:B:305:ILE:HD13	1.99	0.43
1:A:206:PRO:HD3	1:A:275:GLY:O	2.18	0.43
1:C:213:GLU:O	1:C:214:PHE:CB	2.63	0.43
1:A:159:LEU:HD23	1:A:159:LEU:HA	1.53	0.43
1:A:190:PRO:C	1:A:202:VAL:HG21	2.43	0.43
1:D:128:SER:O	1:D:285:ILE:HD13	2.19	0.43
2:E:20:DT:H5'	2:E:20:DT:H6	1.84	0.43
1:B:172:ILE:HD12	1:B:265:TYR:HD2	1.83	0.43
1:B:264:LEU:HD12	1:B:264:LEU:HA	1.79	0.43
1:C:143:GLN:HB3	1:C:173:GLN:OE1	2.18	0.43
1:A:158:GLU:OE1	1:A:158:GLU:N	2.44	0.43
1:B:287:THR:O	1:B:287:THR:HG23	2.17	0.43
1:D:239:ASP:OD1	1:D:241:ILE:HG13	2.19	0.43
1:C:221:PRO:HA	1:C:222:PRO:HD2	1.91	0.43
1:C:285:ILE:HG12	1:C:300:CYS:HG	1.80	0.43
1:D:236:TYR:CD2	1:D:247:VAL:HB	2.53	0.43
2:F:16:DC:H2''	2:F:17:DA:C8	2.54	0.43
1:B:308:CYS:O	1:B:309:PRO:C	2.61	0.43
1:C:227:ARG:NH1	1:C:268:MET:HE2	2.34	0.43
1:D:187:ARG:O	1:D:286:VAL:HA	2.19	0.43
2:E:4:DC:H2''	2:E:5:DA:O5'	2.18	0.43
1:A:142:PHE:HE2	1:A:155:TYR:CE1	2.37	0.42
1:B:163:TYR:N	1:B:163:TYR:CD1	2.87	0.42
1:C:127:PRO:HA	4:C:377:HOH:O	2.18	0.42
1:C:130:THR:O	1:C:298:ARG:NH1	2.52	0.42
1:C:142:PHE:CE1	1:C:172:ILE:HG12	2.55	0.42
1:A:174:ILE:HG13	1:A:186:ILE:HD13	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:180:PRO:HG2	1:B:251:TYR:CD2	2.55	0.42
1:D:151:ALA:HB1	4:D:392:HOH:O	2.17	0.42
1:A:276:GLY:HA2	4:A:372:HOH:O	2.18	0.42
1:C:161:LYS:HE3	1:C:302:GLU:OE2	2.18	0.42
2:E:15:DA:C1'	2:E:16:DC:H5''	2.49	0.42
1:A:154:THR:HG22	1:A:309:PRO:HB2	2.01	0.42
1:A:272:SER:HA	1:A:279:ARG:H	1.84	0.42
1:B:164:CYS:O	1:B:305:ILE:HA	2.19	0.42
1:B:264:LEU:HB2	4:B:358:HOH:O	2.20	0.42
1:C:195:ALA:HA	1:C:198:VAL:CG1	2.50	0.42
1:C:314:LYS:HD2	1:C:314:LYS:O	2.20	0.42
2:F:12:DT:H2''	2:F:13:DA:OP2	2.19	0.42
1:A:145:SER:O	1:A:147:THR:N	2.53	0.42
1:A:318:ASP:C	1:A:320:ILE:N	2.78	0.42
1:D:284:ILE:HB	1:D:301:PHE:CZ	2.55	0.42
2:F:12:DT:C2'	2:F:13:DA:OP2	2.67	0.42
1:B:291:ARG:HH11	1:B:291:ARG:HB2	1.84	0.42
1:C:140:VAL:HA	1:C:173:GLN:O	2.20	0.42
1:C:291:ARG:NH1	1:C:291:ARG:HB3	2.35	0.42
1:B:142:PHE:CE2	1:B:155:TYR:CD1	3.09	0.41
1:A:152:THR:HG22	1:A:153:TRP:CD1	2.56	0.41
1:A:187:ARG:O	1:A:286:VAL:HA	2.20	0.41
1:A:221:PRO:HA	1:A:222:PRO:HD2	1.69	0.41
1:D:159:LEU:O	1:D:160:LYS:C	2.63	0.41
1:D:166:ILE:O	1:D:168:LYS:N	2.52	0.41
1:B:165:GLN:HA	1:B:306:CYS:O	2.20	0.41
1:C:204:ARG:HD3	1:C:224:HIS:O	2.20	0.41
1:A:132:TYR:O	1:A:297:GLY:HA2	2.20	0.41
1:A:204:ARG:CG	1:A:269:CYS:SG	3.08	0.41
1:C:162:LEU:HD22	1:C:162:LEU:C	2.46	0.41
1:C:224:HIS:CD2	1:C:245:GLN:CD	2.99	0.41
1:B:195:ALA:O	1:B:198:VAL:HG22	2.21	0.41
1:A:295:VAL:HG21	1:A:298:ARG:NH2	2.35	0.41
1:D:270:ASN:O	1:D:277:MET:HE3	2.21	0.41
2:E:18:DT:H1'	2:E:19:DG:C8	2.55	0.41
1:B:161:LYS:HD3	1:B:163:TYR:OH	2.20	0.41
1:C:259:GLU:HG3	1:C:260:PHE:CD2	2.56	0.41
1:C:291:ARG:HB3	1:C:291:ARG:CZ	2.51	0.41
2:E:8:DT:C1'	2:E:9:DT:H5''	2.51	0.41
1:A:274:VAL:O	1:A:274:VAL:CG1	2.66	0.41
1:B:227:ARG:HA	1:B:236:TYR:HE2	1.86	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:142:PHE:HE1	1:D:172:ILE:HG12	1.78	0.41
1:A:156:SER:HB3	1:A:159:LEU:HB2	2.02	0.41
1:A:314:LYS:C	1:A:314:LYS:CD	2.94	0.41
1:B:166:ILE:HG12	1:B:167:ALA:N	2.36	0.40
1:C:204:ARG:CG	1:C:269:CYS:SG	3.09	0.40
1:C:308:CYS:HB3	4:C:96:HOH:O	2.20	0.40
1:A:237:VAL:CG2	1:A:238:GLU:N	2.85	0.40
1:C:128:SER:O	1:C:298:ARG:HD3	2.21	0.40
1:C:239:ASP:O	1:C:243:GLY:N	2.53	0.40
1:C:312:ASP:O	1:C:315:ALA:HB3	2.21	0.40
1:B:143:GLN:CB	4:B:380:HOH:O	2.68	0.40
1:B:197:HIS:CG	1:B:280:ARG:HD2	2.57	0.40
1:D:152:THR:O	1:D:153:TRP:HB3	2.21	0.40
1:A:239:ASP:OD1	1:A:242:THR:OG1	2.34	0.40
1:A:283:LEU:HB3	1:A:302:GLU:HA	2.02	0.40
1:B:320:ILE:O	1:B:322:LYS:N	2.51	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	193/203 (95%)	163 (84%)	20 (10%)	10 (5%)	1	1
1	B	189/203 (93%)	174 (92%)	9 (5%)	6 (3%)	3	4
1	C	194/203 (96%)	172 (89%)	15 (8%)	7 (4%)	2	3
1	D	188/203 (93%)	174 (93%)	10 (5%)	4 (2%)	5	9
All	All	764/812 (94%)	683 (89%)	54 (7%)	27 (4%)	3	4

All (27) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	149	LYS
1	A	216	GLU
1	B	151	ALA
1	B	182	GLN
1	B	216	GLU
1	C	146	SER
1	C	147	THR
1	C	148	ALA
1	C	181	PRO
1	D	181	PRO
1	D	216	GLU
1	A	146	SER
1	A	181	PRO
1	B	125	ALA
1	B	181	PRO
1	C	198	VAL
1	A	182	GLN
1	A	198	VAL
1	A	219	ILE
1	C	216	GLU
1	D	167	ALA
1	D	182	GLN
1	A	147	THR
1	A	211	SER
1	A	222	PRO
1	B	167	ALA
1	C	127	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	
1	A	166/179 (93%)	135 (81%)	31 (19%)	1	3
1	B	165/179 (92%)	137 (83%)	28 (17%)	2	4
1	C	168/179 (94%)	143 (85%)	25 (15%)	3	6
1	D	163/179 (91%)	139 (85%)	24 (15%)	3	6

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	662/716 (92%)	554 (84%)	108 (16%)	2 4

All (108) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	144	GLN
1	A	145	SER
1	A	152	THR
1	A	162	LEU
1	A	169	THR
1	A	172	ILE
1	A	173	GLN
1	A	177	MET
1	A	191	VAL
1	A	202	VAL
1	A	203	LYS
1	A	204	ARG
1	A	212	ARG
1	A	218	GLN
1	A	223	SER
1	A	225	LEU
1	A	228	VAL
1	A	235	GLN
1	A	242	THR
1	A	264	LEU
1	A	269	CYS
1	A	271	SER
1	A	274	VAL
1	A	278	ASN
1	A	279	ARG
1	A	283	LEU
1	A	286	VAL
1	A	287	THR
1	A	291	ARG
1	A	296	LEU
1	A	305	ILE
1	B	130	THR
1	B	152	THR
1	B	160	LYS
1	B	162	LEU
1	B	166	ILE
1	B	169	THR

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Mol	Chain	Res	Type
1	B	170	CYS
1	B	177	MET
1	B	194	LYS
1	B	196	GLU
1	B	198	VAL
1	B	215	ASN
1	B	228	VAL
1	B	232	SER
1	B	237	VAL
1	B	241	ILE
1	B	248	LEU
1	B	249	VAL
1	B	262	THR
1	B	264	LEU
1	B	269	CYS
1	B	278	ASN
1	B	279	ARG
1	B	286	VAL
1	B	290	THR
1	B	295	VAL
1	B	296	LEU
1	B	319	SER
1	C	130	THR
1	C	131	ASP
1	C	143	GLN
1	C	147	THR
1	C	152	THR
1	C	160	LYS
1	C	162	LEU
1	C	169	THR
1	C	177	MET
1	C	182	GLN
1	C	201	VAL
1	C	204	ARG
1	C	216	GLU
1	C	219	ILE
1	C	241	ILE
1	C	249	VAL
1	C	264	LEU
1	C	273	CYS
1	C	278	ASN
1	C	286	VAL

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Mol	Chain	Res	Type
1	C	291	ARG
1	C	295	VAL
1	C	296	LEU
1	C	300	CYS
1	C	314	LYS
1	D	124	MET
1	D	130	THR
1	D	143	GLN
1	D	152	THR
1	D	162	LEU
1	D	170	CYS
1	D	177	MET
1	D	182	GLN
1	D	196	GLU
1	D	198	VAL
1	D	228	VAL
1	D	232	SER
1	D	237	VAL
1	D	241	ILE
1	D	248	LEU
1	D	249	VAL
1	D	258	THR
1	D	264	LEU
1	D	278	ASN
1	D	279	ARG
1	D	286	VAL
1	D	295	VAL
1	D	296	LEU
1	D	320	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (24) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	218	GLN
1	A	235	GLN
1	A	245	GLN
1	A	270	ASN
1	B	143	GLN
1	B	173	GLN
1	B	197	HIS
1	B	215	ASN
1	B	218	GLN

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Mol	Chain	Res	Type
1	B	224	HIS
1	B	266	ASN
1	B	270	ASN
1	C	143	GLN
1	C	173	GLN
1	C	182	GLN
1	C	224	HIS
1	C	266	ASN
1	D	182	GLN
1	D	197	HIS
1	D	207	ASN
1	D	218	GLN
1	D	224	HIS
1	D	266	ASN
1	D	278	ASN

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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 4 ligands modelled in this entry, 4 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

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, 95th 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 ²)	Q<0.9
1	A	195/203 (96%)	0.03	5 (2%) 57 52	34, 63, 106, 158	0
1	B	193/203 (95%)	-0.07	2 (1%) 79 76	36, 57, 103, 162	0
1	C	196/203 (96%)	0.31	8 (4%) 41 37	41, 73, 114, 146	0
1	D	192/203 (94%)	0.08	2 (1%) 79 76	37, 67, 113, 138	0
2	E	22/22 (100%)	-0.73	0 100 100	40, 61, 76, 79	0
2	F	22/22 (100%)	-0.73	0 100 100	48, 59, 71, 72	0
All	All	820/856 (95%)	0.04	17 (2%) 63 59	34, 65, 112, 162	0

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	320	ILE	4.5
1	B	151	ALA	4.3
1	C	149	LYS	3.9
1	A	216	GLU	3.8
1	D	151	ALA	3.3
1	C	147	THR	3.1
1	C	300	CYS	2.9
1	A	148	ALA	2.9
1	D	212	ARG	2.8
1	C	125	ALA	2.8
1	A	320	ILE	2.7
1	C	148	ALA	2.7
1	C	241	ILE	2.2
1	A	147	THR	2.1
1	A	149	LYS	2.0
1	B	124	MET	2.0
1	C	126	SER	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	ZN	C	901	1/1	0.99	0.06	79,79,79,79	0
3	ZN	B	901	1/1	1.00	0.03	51,51,51,51	0
3	ZN	A	901	1/1	1.00	0.04	54,54,54,54	0
3	ZN	D	901	1/1	1.00	0.02	59,59,59,59	0

6.5 Other polymers [i](#)

There are no such residues in this entry.