



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 03:29 PM UTC

PDB ID : 2Z9O / pdb_00002z9o
Title : Crystal structure of the dimeric form of RepE in complex with the repE operator DNA
Authors : Nakamura, A.; Wada, C.; Miki, K.
Deposited on : 2007-09-21
Resolution : 3.14 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

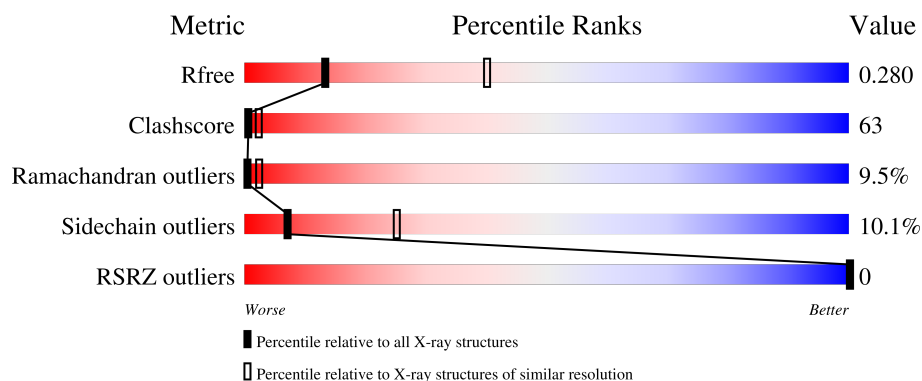
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

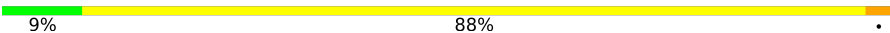
The reported resolution of this entry is 3.14 Å.

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	2351 (3.18-3.10)
Clashscore	190562	2452 (3.18-3.10)
Ramachandran outliers	187476	2324 (3.18-3.10)
Sidechain outliers	187428	2324 (3.18-3.10)
RSRZ outliers	180081	2351 (3.18-3.10)

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 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	C	33	
2	D	33	
3	A	266	
3	B	266	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4926 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 (33-MER).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	C	33	Total	C	N	O	P	0	0	0
			669	323	118	196	32			

- Molecule 2 is a DNA chain called DNA (33-MER).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	33	Total	C	N	O	P	0	0	0
			678	326	124	196	32			

- Molecule 3 is a protein called Replication initiation protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	A	222	Total	C	N	O	S	0	0	0
			1796	1151	312	325	8			
3	B	218	Total	C	N	O	S	0	0	0
			1783	1142	310	324	7			

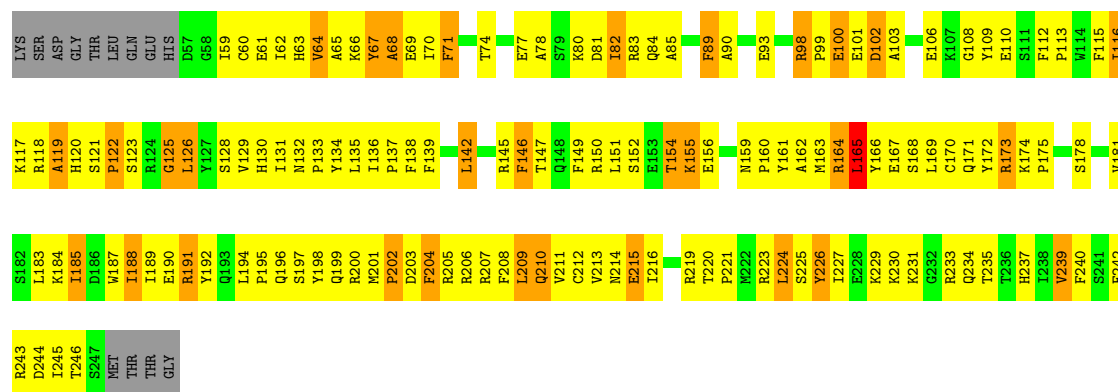
There are 32 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-14	MET	-	expression tag	UNP P03856
A	-13	ARG	-	expression tag	UNP P03856
A	-12	GLY	-	expression tag	UNP P03856
A	-11	SER	-	expression tag	UNP P03856
A	-10	HIS	-	expression tag	UNP P03856
A	-9	HIS	-	expression tag	UNP P03856
A	-8	HIS	-	expression tag	UNP P03856
A	-7	HIS	-	expression tag	UNP P03856
A	-6	HIS	-	expression tag	UNP P03856
A	-5	HIS	-	expression tag	UNP P03856
A	-4	GLY	-	expression tag	UNP P03856
A	-3	SER	-	expression tag	UNP P03856

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	ILE	-	expression tag	UNP P03856
A	-1	GLU	-	expression tag	UNP P03856
A	0	GLY	-	expression tag	UNP P03856
A	1	ARG	-	expression tag	UNP P03856
B	-14	MET	-	expression tag	UNP P03856
B	-13	ARG	-	expression tag	UNP P03856
B	-12	GLY	-	expression tag	UNP P03856
B	-11	SER	-	expression tag	UNP P03856
B	-10	HIS	-	expression tag	UNP P03856
B	-9	HIS	-	expression tag	UNP P03856
B	-8	HIS	-	expression tag	UNP P03856
B	-7	HIS	-	expression tag	UNP P03856
B	-6	HIS	-	expression tag	UNP P03856
B	-5	HIS	-	expression tag	UNP P03856
B	-4	GLY	-	expression tag	UNP P03856
B	-3	SER	-	expression tag	UNP P03856
B	-2	ILE	-	expression tag	UNP P03856
B	-1	GLU	-	expression tag	UNP P03856
B	0	GLY	-	expression tag	UNP P03856
B	1	ARG	-	expression tag	UNP P03856



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	60.73Å 99.32Å 95.00Å 90.00° 108.55° 90.00°	Depositor
Resolution (Å)	45.03 – 3.14 45.03 – 3.14	Depositor EDS
% Data completeness (in resolution range)	87.8 (45.03-3.14) 87.8 (45.03-3.14)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.01 (at 3.12Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.266 , 0.313 0.247 , 0.280	Depositor DCC
R_{free} test set	850 reflections (5.15%)	wwPDB-VP
Wilson B-factor (Å ²)	51.2	Xtriage
Anisotropy	1.080	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 43.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.42$, $\langle L^2 \rangle = 0.25$	Xtriage
Estimated twinning fraction	0.088 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4926	wwPDB-VP
Average B, all atoms (Å ²)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.43% 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

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	0.37	0/749	0.74	0/1153
2	D	0.35	0/761	0.70	0/1174
3	A	0.54	0/1839	1.06	17/2480 (0.7%)
3	B	0.54	0/1826	1.01	10/2462 (0.4%)
All	All	0.50	0/5175	0.95	27/7269 (0.4%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	C	0	1

There are no bond length outliers.

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	205	ARG	N-CA-C	7.12	118.73	110.97
3	A	37	ARG	N-CA-C	-7.10	103.14	111.03
3	B	116	ILE	N-CA-C	-6.91	105.11	111.67
3	A	44	ASP	N-CA-C	-6.87	103.87	111.36
3	B	146	PHE	N-CA-C	-6.74	104.02	111.36
3	B	239	VAL	N-CA-C	6.58	116.09	106.55
3	A	135	LEU	N-CA-C	6.46	119.63	111.69
3	A	224	LEU	N-CA-C	6.08	119.03	108.76
3	A	236	THR	N-CA-C	6.04	120.75	112.04
3	A	75	SER	N-CA-C	-5.98	104.69	112.23
3	B	142	LEU	N-CA-C	-5.84	105.05	111.82
3	A	162	ALA	N-CA-C	-5.71	104.74	110.97
3	B	71	PHE	N-CA-C	5.64	119.79	113.02
3	A	73	LEU	N-CA-C	5.38	117.27	108.76

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	27	ALA	N-CA-C	-5.38	106.40	113.12
3	A	119	ALA	N-CA-C	5.38	116.69	108.52
3	A	233	ARG	N-CA-C	-5.36	107.39	114.04
3	A	201	MET	N-CA-C	5.36	121.65	109.81
3	A	199	GLN	N-CA-C	-5.29	106.34	112.89
3	A	31	LEU	N-CA-C	5.21	118.18	110.46
3	B	119	ALA	N-CA-C	5.12	116.30	108.42
3	B	201	MET	N-CA-C	5.11	120.62	113.57
3	A	192	TYR	N-CA-C	-5.08	105.44	113.02
3	B	102	ASP	N-CA-C	5.08	117.36	110.35
3	B	155	LYS	N-CA-C	-5.08	101.82	109.60
3	A	176	ASP	N-CA-C	-5.06	106.02	113.61
3	B	224	LEU	N-CA-C	5.06	117.03	109.59

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	1	DT	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	C	669	0	376	71	0
2	D	678	0	376	67	0
3	A	1796	0	1742	236	0
3	B	1783	0	1735	240	0
All	All	4926	0	4229	573	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 63.

All (573) 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)
3:B:200:ARG:HG2	3:B:202:PRO:HD2	1.15	1.10
2:D:39:DG:H2''	2:D:40:DT:H5''	1.25	1.10
2:D:41:DG:H2''	2:D:42:DA:H5'	1.38	1.04
1:C:6:DG:C2'	1:C:7:DT:H5''	1.86	1.04
2:D:60:DA:H2''	2:D:61:DC:H5'	1.40	1.04
3:B:38:MET:HE1	3:B:82:ILE:HG23	1.36	1.04
3:A:119:ALA:HB2	3:B:110:GLU:HA	1.42	0.99
2:D:39:DG:C2'	2:D:40:DT:H5''	1.93	0.99
1:C:6:DG:H2''	1:C:7:DT:H5''	1.43	0.98
3:B:164:ARG:HG2	3:B:164:ARG:HH11	1.25	0.98
1:C:24:DG:H2''	1:C:25:DT:H5'	1.45	0.98
3:B:63:HIS:HB2	3:B:66:LYS:HG2	1.45	0.97
3:A:42:PHE:HD1	3:A:131:ILE:HD11	1.30	0.96
3:A:244:ASP:OD2	3:A:246:THR:HG23	1.66	0.95
1:C:31:DT:H2''	1:C:32:DT:H5'	1.49	0.94
1:C:8:DG:H2''	1:C:9:DA:H5'	1.48	0.94
3:A:122:PRO:HG2	3:B:102:ASP:HA	1.49	0.93
2:D:60:DA:H62	3:B:200:ARG:HH12	1.12	0.92
3:B:223:ARG:HD3	3:B:245:ILE:HG13	1.51	0.92
3:A:132:ASN:HD22	3:A:132:ASN:C	1.77	0.90
1:C:1:DT:H2''	1:C:2:DT:H5'	1.54	0.89
1:C:21:DC:H2''	1:C:22:DT:H5''	1.54	0.89
3:B:206:ARG:O	3:B:207:ARG:HD3	1.73	0.88
2:D:50:DT:H2''	2:D:51:DT:H5'	1.55	0.88
3:A:116:ILE:HG12	3:B:116:ILE:HD12	1.56	0.88
1:C:6:DG:H2''	1:C:7:DT:C5'	2.03	0.87
3:B:98:ARG:HG2	3:B:98:ARG:HH11	1.40	0.86
3:B:37:ARG:HD2	3:B:71:PHE:CE2	2.10	0.86
1:C:22:DT:H2''	1:C:23:DT:H5'	1.56	0.85
3:A:42:PHE:O	3:A:46:ILE:HG12	1.75	0.85
3:A:132:ASN:HD22	3:A:133:PRO:N	1.74	0.85
3:A:160:PRO:O	3:A:164:ARG:HG3	1.76	0.84
1:C:4:DG:H5'	3:B:233:ARG:HB3	1.59	0.84
3:B:198:TYR:HE1	3:B:207:ARG:HB2	1.41	0.83
3:A:42:PHE:H	3:A:42:PHE:HD2	1.25	0.83
2:D:64:DT:H2''	2:D:65:DA:C8	2.14	0.83
1:C:29:DA:H2''	1:C:30:DC:H5''	1.60	0.83
2:D:52:DA:H1'	2:D:53:DG:H5''	1.60	0.82
1:C:7:DT:H4'	1:C:7:DT:OP1	1.79	0.82
3:A:164:ARG:HH11	3:A:164:ARG:HG2	1.45	0.80
3:A:145:ARG:HG2	3:A:145:ARG:HH11	1.46	0.80
3:A:42:PHE:CD1	3:A:131:ILE:HD11	2.15	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:99:PRO:HG2	3:A:100:GLU:H	1.45	0.80
3:A:37:ARG:HB3	3:A:71:PHE:CE1	2.17	0.79
3:B:160:PRO:HA	3:B:163:MET:HG2	1.64	0.79
1:C:21:DC:H2'	1:C:22:DT:H71	1.63	0.79
3:A:213:VAL:HG13	3:A:224:LEU:HB3	1.63	0.79
3:B:116:ILE:HG12	3:B:132:ASN:HA	1.64	0.79
3:A:32:SER:OG	3:A:35:GLN:HG3	1.83	0.79
3:B:26:GLU:HB2	3:B:163:MET:SD	2.23	0.79
3:A:119:ALA:CB	3:B:110:GLU:HA	2.13	0.78
3:A:62:ILE:HG21	3:A:82:ILE:HD12	1.64	0.78
3:B:159:ASN:HD22	3:B:160:PRO:HD2	1.48	0.78
3:A:116:ILE:HD11	3:A:133:PRO:HD2	1.65	0.78
3:A:73:LEU:HD21	3:A:77:GLU:OE1	1.84	0.77
1:C:4:DG:C5'	3:B:233:ARG:HB3	2.14	0.77
3:B:187:TRP:CZ2	3:B:191:ARG:HG3	2.19	0.77
2:D:45:DA:H2''	2:D:46:DG:H5'	1.66	0.77
3:A:196:GLN:HG3	3:A:197:SER:N	1.99	0.77
2:D:41:DG:H2''	2:D:42:DA:C5'	2.15	0.76
3:A:31:LEU:HD22	3:A:35:GLN:NE2	2.00	0.76
3:A:73:LEU:HD22	3:A:77:GLU:HB2	1.67	0.76
3:B:154:THR:HG23	3:B:220:THR:HG22	1.66	0.75
3:B:173:ARG:O	3:B:174:LYS:HG3	1.86	0.75
2:D:60:DA:H2''	2:D:61:DC:C5'	2.16	0.75
2:D:60:DA:N6	3:B:200:ARG:HH12	1.84	0.75
3:A:116:ILE:HG13	3:A:132:ASN:HA	1.69	0.75
2:D:42:DA:H2''	2:D:43:DC:H5'	1.69	0.74
2:D:58:DT:H2''	2:D:59:DC:H5''	1.68	0.74
3:B:98:ARG:HB3	3:B:101:GLU:HB2	1.70	0.74
3:B:42:PHE:HE2	3:B:62:ILE:HD11	1.53	0.74
3:B:100:GLU:HG3	3:B:101:GLU:N	2.01	0.74
1:C:29:DA:C2'	1:C:30:DC:H5''	2.18	0.74
3:A:249:THR:O	3:A:249:THR:HG22	1.87	0.73
1:C:17:DA:H1'	1:C:18:DA:H5''	1.70	0.73
3:A:68:ALA:HB1	3:A:73:LEU:O	1.88	0.73
3:A:122:PRO:CG	3:B:102:ASP:HA	2.19	0.73
1:C:1:DT:H3	2:D:66:DA:H2	1.37	0.73
3:B:200:ARG:HG2	3:B:202:PRO:CD	2.08	0.73
3:B:164:ARG:HG2	3:B:164:ARG:NH1	2.02	0.72
3:B:74:THR:HG23	3:B:77:GLU:CD	2.13	0.72
3:B:172:TYR:HD1	3:B:181:VAL:HG21	1.54	0.72
3:B:61:GLU:OE1	3:B:126:LEU:HD11	1.90	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:34:ASP:HA	3:A:37:ARG:HG3	1.72	0.72
3:A:150:ARG:HH22	3:A:246:THR:HG21	1.53	0.72
3:B:36:LYS:HE2	3:B:154:THR:O	1.90	0.71
3:B:196:GLN:CD	3:B:196:GLN:H	1.97	0.71
3:A:150:ARG:HH22	3:A:246:THR:CG2	2.02	0.71
1:C:12:DA:C2'	1:C:13:DT:H5''	2.21	0.71
3:A:223:ARG:HB3	3:A:243:ARG:NH1	2.06	0.71
3:A:228:GLU:HB3	3:A:235:THR:HG21	1.73	0.71
2:D:55:DT:H2''	2:D:56:DT:C5'	2.21	0.71
3:B:100:GLU:HG3	3:B:101:GLU:H	1.54	0.71
3:B:229:LYS:HG3	3:B:229:LYS:O	1.91	0.70
1:C:27:DA:H2''	1:C:28:DC:H5''	1.71	0.70
3:B:64:VAL:HG22	3:B:125:GLY:O	1.92	0.70
3:A:32:SER:HA	3:A:156:GLU:OE2	1.91	0.70
3:A:119:ALA:HB1	3:B:109:TYR:O	1.90	0.70
2:D:39:DG:H2''	2:D:40:DT:C5'	2.15	0.69
2:D:54:DA:H2''	2:D:55:DT:H5'	1.73	0.69
3:A:189:ILE:HD13	3:A:199:GLN:HG2	1.74	0.69
3:B:184:LYS:HD2	3:B:237:HIS:CD2	2.28	0.69
2:D:49:DT:H2''	2:D:50:DT:H5''	1.75	0.69
3:A:168:SER:O	3:A:171:GLN:HB3	1.93	0.69
3:B:102:ASP:HB2	3:B:106:GLU:O	1.93	0.69
3:B:227:ILE:HB	3:B:239:VAL:HB	1.74	0.69
3:A:100:GLU:HG2	3:A:101:GLU:N	2.07	0.69
1:C:12:DA:H2''	1:C:13:DT:C5'	2.23	0.69
3:A:178:SER:HA	3:A:243:ARG:HA	1.75	0.69
3:B:21:SER:O	3:B:22:ASN:HB3	1.93	0.69
3:B:200:ARG:CG	3:B:202:PRO:HD2	2.09	0.69
3:A:243:ARG:HG3	3:A:248:MET:SD	2.33	0.69
2:D:39:DG:C3'	2:D:40:DT:H5''	2.23	0.68
1:C:6:DG:H2'	1:C:7:DT:H5''	1.75	0.68
2:D:59:DC:H2''	2:D:60:DA:C8	2.27	0.68
1:C:31:DT:H2''	1:C:32:DT:C5'	2.22	0.68
2:D:54:DA:H2''	2:D:55:DT:C5'	2.24	0.68
3:A:145:ARG:HH11	3:A:145:ARG:CG	2.06	0.68
2:D:58:DT:C2'	2:D:59:DC:H5''	2.23	0.68
1:C:12:DA:H2''	1:C:13:DT:H5''	1.76	0.68
3:B:160:PRO:HA	3:B:163:MET:CG	2.24	0.67
3:B:116:ILE:HD11	3:B:133:PRO:HD2	1.77	0.67
3:B:160:PRO:HB2	3:B:164:ARG:CZ	2.25	0.67
3:B:198:TYR:CE1	3:B:207:ARG:HB2	2.27	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:102:ASP:HB3	3:A:108:GLY:H	1.59	0.67
3:B:98:ARG:HG2	3:B:98:ARG:NH1	2.07	0.67
3:B:230:LYS:HA	3:B:235:THR:HA	1.76	0.67
3:A:172:TYR:CB	3:A:181:VAL:HG21	2.25	0.67
2:D:58:DT:H2''	2:D:59:DC:C5'	2.25	0.67
3:B:160:PRO:O	3:B:164:ARG:HG3	1.95	0.67
1:C:21:DC:H2''	1:C:22:DT:C5'	2.25	0.66
3:B:102:ASP:HB3	3:B:108:GLY:N	2.09	0.66
3:B:160:PRO:HB2	3:B:164:ARG:NH2	2.10	0.66
3:A:116:ILE:CD1	3:A:133:PRO:HD2	2.25	0.66
2:D:61:DC:H2''	2:D:62:DA:C8	2.30	0.66
3:B:110:GLU:HG2	3:B:112:PHE:CZ	2.31	0.66
3:A:245:ILE:HD13	3:A:245:ILE:O	1.96	0.66
3:B:164:ARG:HH11	3:B:164:ARG:CG	2.07	0.65
2:D:57:DG:H1'	2:D:58:DT:H5''	1.79	0.65
3:A:208:PHE:HD1	3:A:209:LEU:N	1.94	0.65
3:B:133:PRO:HA	3:B:136:ILE:CD1	2.26	0.65
3:B:200:ARG:HE	3:B:202:PRO:HG2	1.62	0.65
3:A:97:TYR:O	3:A:98:ARG:HD2	1.96	0.65
3:A:136:ILE:N	3:A:137:PRO:HD2	2.11	0.65
1:C:6:DG:H2''	1:C:7:DT:C4'	2.26	0.64
1:C:14:DC:H2''	1:C:15:DT:O5'	1.97	0.64
1:C:12:DA:H1'	1:C:13:DT:H5''	1.78	0.64
3:A:42:PHE:HD2	3:A:42:PHE:N	1.93	0.64
3:A:200:ARG:HH12	3:A:202:PRO:HB2	1.63	0.64
1:C:24:DG:H2''	1:C:25:DT:C5'	2.25	0.63
1:C:26:DC:H2''	1:C:27:DA:C8	2.32	0.63
3:B:230:LYS:HG3	3:B:234:GLN:C	2.22	0.63
2:D:55:DT:H2''	2:D:56:DT:H5'	1.78	0.63
3:B:209:LEU:O	3:B:210:GLN:C	2.41	0.63
3:A:174:LYS:O	3:A:176:ASP:N	2.31	0.63
3:B:126:LEU:C	3:B:126:LEU:HD23	2.23	0.63
3:A:122:PRO:HG2	3:B:102:ASP:CA	2.26	0.63
3:B:212:CYS:O	3:B:216:ILE:HG13	1.99	0.63
3:A:27:ALA:O	3:A:30:SER:HB3	1.99	0.63
3:A:174:LYS:C	3:A:176:ASP:H	2.06	0.63
3:A:249:THR:O	3:A:249:THR:CG2	2.47	0.63
2:D:38:DT:H2''	2:D:39:DG:OP2	1.98	0.62
3:A:200:ARG:NH1	3:A:202:PRO:HB2	2.14	0.62
3:A:110:GLU:CD	3:B:117:LYS:HE3	2.23	0.62
3:B:170:CYS:O	3:B:173:ARG:HG3	1.98	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:60:DA:H62	3:B:200:ARG:NH1	1.91	0.62
3:A:150:ARG:NH2	3:A:246:THR:HG21	2.12	0.62
3:A:116:ILE:O	3:B:113:PRO:HD2	2.00	0.62
3:A:123:SER:HB3	3:B:101:GLU:O	2.00	0.62
3:A:169:LEU:HD13	3:A:224:LEU:HD11	1.81	0.62
3:A:172:TYR:CG	3:A:181:VAL:HG21	2.34	0.62
3:A:223:ARG:NE	3:A:245:ILE:HG13	2.13	0.62
3:B:20:GLN:O	3:B:145:ARG:NH1	2.33	0.62
3:B:66:LYS:O	3:B:70:ILE:HG12	1.99	0.62
3:A:93:GLU:OE2	3:B:118:ARG:NH2	2.33	0.62
3:A:116:ILE:HG12	3:B:116:ILE:CD1	2.29	0.62
3:A:165:LEU:CD2	3:A:169:LEU:HG	2.30	0.61
3:A:224:LEU:CD1	3:A:242:PHE:HB3	2.29	0.61
3:A:165:LEU:HA	3:A:192:TYR:CZ	2.35	0.61
3:B:42:PHE:CE2	3:B:62:ILE:HD11	2.34	0.61
1:C:25:DT:H2''	1:C:26:DC:H5''	1.81	0.61
3:A:194:LEU:HD21	3:A:208:PHE:CD2	2.34	0.61
3:A:208:PHE:CD1	3:A:209:LEU:N	2.68	0.61
3:A:224:LEU:HD12	3:A:242:PHE:HB3	1.81	0.61
3:B:128:SER:O	3:B:129:VAL:HG23	2.00	0.61
3:B:146:PHE:HD2	3:B:152:SER:HB2	1.65	0.61
3:B:172:TYR:HD1	3:B:181:VAL:CG2	2.13	0.61
3:B:195:PRO:O	3:B:198:TYR:HB2	2.00	0.61
3:B:195:PRO:HD2	3:B:198:TYR:CD2	2.35	0.61
2:D:44:DA:H2''	2:D:45:DA:H5'	1.83	0.61
3:B:59:ILE:HG22	3:B:60:CYS:N	2.13	0.61
1:C:4:DG:H5''	3:B:233:ARG:C	2.26	0.60
3:B:24:LEU:O	3:B:24:LEU:HD23	2.01	0.60
3:A:184:LYS:HA	3:A:237:HIS:CD2	2.36	0.60
1:C:16:DA:H2''	1:C:17:DA:OP2	2.01	0.60
2:D:45:DA:H1'	2:D:46:DG:H5''	1.83	0.60
3:B:146:PHE:HA	3:B:151:LEU:H	1.66	0.60
3:B:183:LEU:HD13	3:B:187:TRP:HZ3	1.64	0.60
1:C:25:DT:C2'	1:C:26:DC:H5''	2.32	0.60
2:D:50:DT:C2'	2:D:51:DT:H5'	2.31	0.60
3:B:78:ALA:O	3:B:82:ILE:HG12	2.02	0.60
1:C:4:DG:OP1	3:B:234:GLN:HB2	2.01	0.59
3:A:140:ILE:O	3:A:143:GLN:HG2	2.01	0.59
3:A:123:SER:CB	3:B:101:GLU:O	2.51	0.59
3:A:196:GLN:HG3	3:A:197:SER:H	1.68	0.59
2:D:57:DG:H2''	2:D:58:DT:H5'	1.84	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:67:TYR:O	3:A:70:ILE:HG22	2.03	0.59
3:B:67:TYR:O	3:B:70:ILE:N	2.34	0.59
1:C:24:DG:H1'	1:C:25:DT:H5''	1.85	0.59
3:B:160:PRO:O	3:B:163:MET:HG2	2.03	0.59
3:B:192:TYR:HB2	3:B:194:LEU:HD13	1.84	0.58
3:A:164:ARG:HG2	3:A:164:ARG:NH1	2.14	0.58
3:B:22:ASN:O	3:B:25:THR:HB	2.02	0.58
3:B:63:HIS:HB2	3:B:66:LYS:CG	2.27	0.58
3:B:37:ARG:HD2	3:B:71:PHE:HE2	1.67	0.58
3:B:37:ARG:NH1	3:B:71:PHE:CD2	2.72	0.58
2:D:35:DA:H1'	2:D:36:DA:H5''	1.85	0.58
3:A:205:ARG:HG3	3:A:205:ARG:HH11	1.68	0.58
3:B:131:ILE:HD13	3:B:139:PHE:CZ	2.39	0.58
1:C:11:DA:H2''	1:C:12:DA:H5'	1.85	0.57
2:D:45:DA:H2''	2:D:46:DG:C5'	2.33	0.57
3:A:62:ILE:CG2	3:A:82:ILE:HD12	2.34	0.57
3:B:223:ARG:HD3	3:B:245:ILE:CG1	2.30	0.57
3:A:45:GLN:C	3:A:47:ARG:H	2.11	0.57
1:C:25:DT:H2''	1:C:26:DC:C5'	2.35	0.57
3:A:200:ARG:HG2	3:A:202:PRO:HD2	1.87	0.57
1:C:12:DA:C1'	1:C:13:DT:H5''	2.35	0.57
3:A:113:PRO:HD2	3:B:116:ILE:O	2.05	0.57
3:A:81:ASP:O	3:A:82:ILE:C	2.48	0.57
1:C:27:DA:C2'	1:C:28:DC:H5''	2.35	0.56
2:D:62:DA:H1'	2:D:63:DC:H5'	1.86	0.56
3:A:36:LYS:CE	3:A:151:LEU:HD11	2.35	0.56
3:A:37:ARG:O	3:A:38:MET:C	2.48	0.56
1:C:18:DA:H5'	1:C:18:DA:H8	1.70	0.56
3:A:190:GLU:C	3:A:192:TYR:H	2.12	0.56
3:A:34:ASP:CA	3:A:37:ARG:HG3	2.35	0.56
3:A:47:ARG:HH11	3:A:47:ARG:HG2	1.71	0.56
3:A:205:ARG:HG3	3:A:205:ARG:NH1	2.21	0.56
3:B:187:TRP:C	3:B:189:ILE:H	2.11	0.56
1:C:17:DA:C2'	1:C:18:DA:H5''	2.36	0.56
3:B:67:TYR:O	3:B:69:GLU:N	2.39	0.56
3:B:33:ARG:NH2	3:B:34:ASP:OD1	2.39	0.56
3:B:196:GLN:C	3:B:198:TYR:H	2.14	0.56
3:B:187:TRP:CE2	3:B:191:ARG:HG3	2.41	0.56
1:C:29:DA:C3'	1:C:30:DC:H5''	2.36	0.55
3:A:99:PRO:O	3:A:102:ASP:OD1	2.25	0.55
3:B:159:ASN:ND2	3:B:160:PRO:HD2	2.21	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:184:LYS:HG3	3:A:237:HIS:CD2	2.42	0.55
3:A:100:GLU:HG2	3:A:101:GLU:H	1.72	0.55
3:A:223:ARG:O	3:A:242:PHE:HA	2.06	0.55
3:B:160:PRO:HG2	3:B:161:TYR:H	1.72	0.55
3:A:122:PRO:HB2	3:B:103:ALA:HB3	1.89	0.55
3:A:177:GLY:O	3:A:244:ASP:N	2.40	0.55
3:B:32:SER:HA	3:B:156:GLU:OE2	2.06	0.55
3:A:97:TYR:HE2	3:A:107:LYS:HZ2	1.53	0.55
1:C:4:DG:H2''	1:C:5:DT:OP2	2.07	0.54
3:A:100:GLU:CG	3:A:101:GLU:N	2.70	0.54
3:A:117:LYS:HE3	3:B:112:PHE:CE2	2.42	0.54
1:C:17:DA:C1'	1:C:18:DA:H5''	2.36	0.54
3:B:25:THR:C	3:B:27:ALA:H	2.13	0.54
3:B:129:VAL:O	3:B:130:HIS:HD2	1.91	0.54
3:B:160:PRO:CA	3:B:163:MET:HG2	2.35	0.54
2:D:49:DT:C2'	2:D:50:DT:H5''	2.37	0.54
2:D:42:DA:C2'	2:D:43:DC:H5'	2.37	0.54
3:B:102:ASP:HB3	3:B:108:GLY:H	1.70	0.54
3:A:31:LEU:HD22	3:A:35:GLN:HE22	1.71	0.54
3:A:33:ARG:HD2	3:A:153:GLU:HA	1.90	0.54
3:A:174:LYS:HG2	3:A:178:SER:O	2.08	0.54
3:A:179:GLY:O	3:A:241:SER:HA	2.08	0.54
3:A:227:ILE:O	3:A:227:ILE:HG13	2.08	0.54
3:B:32:SER:O	3:B:33:ARG:C	2.51	0.54
3:B:231:LYS:HG3	3:B:231:LYS:O	2.08	0.54
3:A:213:VAL:O	3:A:217:ASN:ND2	2.41	0.54
3:A:223:ARG:HB3	3:A:243:ARG:HH11	1.70	0.54
3:B:229:LYS:HG2	3:B:237:HIS:HB2	1.90	0.54
2:D:63:DC:C2'	2:D:64:DT:H72	2.37	0.53
3:A:133:PRO:O	3:A:134:TYR:C	2.51	0.53
3:B:167:GLU:O	3:B:171:GLN:HG2	2.08	0.53
2:D:63:DC:H2''	2:D:64:DT:H72	1.90	0.53
3:B:196:GLN:CD	3:B:196:GLN:N	2.66	0.53
3:B:243:ARG:NH1	3:B:245:ILE:HD13	2.23	0.53
3:B:67:TYR:HD2	3:B:68:ALA:N	2.07	0.53
3:B:168:SER:O	3:B:169:LEU:C	2.51	0.53
3:B:194:LEU:HD21	3:B:208:PHE:CD2	2.44	0.53
3:A:142:LEU:O	3:A:145:ARG:N	2.41	0.53
1:C:19:DA:H1'	1:C:20:DA:H5'	1.91	0.52
3:B:118:ARG:HG2	3:B:119:ALA:N	2.25	0.52
3:B:155:LYS:HD2	3:B:219:ARG:HD2	1.90	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:229:LYS:CG	3:B:237:HIS:HB2	2.40	0.52
2:D:57:DG:N7	3:B:207:ARG:NH1	2.56	0.52
3:A:25:THR:C	3:A:27:ALA:H	2.17	0.52
1:C:19:DA:H2''	1:C:20:DA:O5'	2.09	0.52
3:B:150:ARG:NH1	3:B:244:ASP:OD1	2.40	0.52
3:B:169:LEU:HB3	3:B:242:PHE:CD2	2.45	0.52
1:C:6:DG:H2''	1:C:7:DT:O4'	2.10	0.52
3:A:142:LEU:O	3:A:145:ARG:HB2	2.10	0.52
3:B:116:ILE:CG1	3:B:132:ASN:HA	2.36	0.52
3:A:22:ASN:OD1	3:A:24:LEU:HB3	2.10	0.52
3:A:165:LEU:HA	3:A:192:TYR:OH	2.10	0.52
3:B:151:LEU:HA	3:B:154:THR:OG1	2.10	0.52
3:B:164:ARG:O	3:B:166:TYR:N	2.43	0.52
3:B:183:LEU:HD13	3:B:187:TRP:CZ3	2.44	0.52
3:A:36:LYS:HE3	3:A:151:LEU:HD11	1.91	0.51
3:A:100:GLU:CG	3:A:101:GLU:H	2.22	0.51
3:B:133:PRO:HA	3:B:136:ILE:HD12	1.90	0.51
3:A:168:SER:HB2	3:A:192:TYR:OH	2.11	0.51
3:B:38:MET:HE2	3:B:62:ILE:HD13	1.93	0.51
1:C:28:DC:H2'	1:C:29:DA:C8	2.46	0.51
3:A:184:LYS:HG3	3:A:237:HIS:NE2	2.25	0.51
2:D:55:DT:H2''	2:D:56:DT:H5''	1.91	0.51
3:B:203:ASP:O	3:B:204:PHE:C	2.53	0.51
3:B:198:TYR:HE1	3:B:207:ARG:CB	2.19	0.51
3:B:229:LYS:CD	3:B:237:HIS:HB2	2.40	0.51
1:C:29:DA:H2''	1:C:30:DC:O4'	2.11	0.51
3:A:129:VAL:HG23	3:A:129:VAL:O	2.11	0.51
3:A:32:SER:O	3:A:35:GLN:HB2	2.11	0.51
3:B:110:GLU:HG2	3:B:112:PHE:HZ	1.73	0.51
3:A:40:TYR:O	3:A:41:LEU:C	2.54	0.50
3:B:35:GLN:HG2	3:B:85:ALA:HB1	1.92	0.50
1:C:16:DA:C2	2:D:52:DA:C2	2.99	0.50
3:A:139:PHE:O	3:A:141:GLY:N	2.44	0.50
1:C:6:DG:C2'	1:C:7:DT:C5'	2.67	0.50
3:B:149:PHE:N	3:B:149:PHE:CD1	2.79	0.50
3:B:164:ARG:NH1	3:B:164:ARG:CG	2.70	0.50
3:B:196:GLN:C	3:B:198:TYR:N	2.68	0.50
3:A:90:ALA:O	3:A:92:LYS:N	2.44	0.50
3:A:42:PHE:N	3:A:42:PHE:CD2	2.65	0.50
3:A:104:GLY:C	3:A:106:GLU:H	2.19	0.50
2:D:43:DC:H2''	2:D:44:DA:C8	2.45	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:102:ASP:HA	3:B:122:PRO:HG2	1.92	0.50
3:A:119:ALA:HA	3:B:109:TYR:CE1	2.47	0.50
3:A:156:GLU:CD	3:A:156:GLU:H	2.20	0.50
3:A:172:TYR:CD1	3:A:181:VAL:HG11	2.47	0.50
3:A:132:ASN:C	3:A:132:ASN:ND2	2.51	0.50
3:B:209:LEU:HD22	3:B:226:TYR:CD2	2.47	0.50
1:C:25:DT:H1'	1:C:26:DC:H5''	1.94	0.50
2:D:60:DA:N6	3:B:200:ARG:HH22	2.10	0.49
3:B:195:PRO:HD2	3:B:198:TYR:HD2	1.77	0.49
3:A:184:LYS:O	3:A:187:TRP:HB3	2.12	0.49
3:A:236:THR:O	3:A:237:HIS:HD2	1.95	0.49
3:B:203:ASP:O	3:B:206:ARG:N	2.45	0.49
2:D:52:DA:C1'	2:D:53:DG:H5''	2.37	0.49
3:A:102:ASP:HB3	3:A:108:GLY:N	2.25	0.49
3:A:120:HIS:ND1	3:A:127:TYR:CE1	2.81	0.49
3:B:74:THR:HG23	3:B:77:GLU:OE2	2.12	0.49
3:A:130:HIS:HD1	3:B:134:TYR:HH	1.59	0.49
3:A:139:PHE:O	3:A:140:ILE:C	2.56	0.49
3:A:212:CYS:O	3:A:213:VAL:C	2.55	0.49
3:B:196:GLN:O	3:B:199:GLN:HG2	2.11	0.49
3:A:146:PHE:HD2	3:A:152:SER:HG	1.59	0.49
3:A:37:ARG:HB3	3:A:71:PHE:HE1	1.73	0.49
3:A:116:ILE:CG1	3:B:116:ILE:HD12	2.36	0.49
3:A:145:ARG:CG	3:A:145:ARG:NH1	2.69	0.49
3:B:59:ILE:CG2	3:B:60:CYS:N	2.75	0.49
3:B:164:ARG:C	3:B:166:TYR:N	2.70	0.49
3:B:59:ILE:HG23	3:B:130:HIS:CD2	2.48	0.48
1:C:4:DG:H5''	3:B:233:ARG:HB3	1.94	0.48
3:A:218:SER:O	3:A:219:ARG:HD3	2.13	0.48
3:A:229:LYS:O	3:A:236:THR:HG23	2.14	0.48
3:B:243:ARG:HH11	3:B:245:ILE:HD13	1.78	0.48
3:A:228:GLU:HB3	3:A:235:THR:CG2	2.42	0.48
3:B:169:LEU:HB3	3:B:242:PHE:CE2	2.47	0.48
3:B:173:ARG:O	3:B:174:LYS:CG	2.58	0.48
1:C:13:DT:H2''	1:C:14:DC:O5'	2.12	0.48
3:A:209:LEU:O	3:A:210:GLN:C	2.55	0.48
3:B:214:ASN:C	3:B:216:ILE:H	2.21	0.48
3:A:112:PHE:CD1	3:A:112:PHE:N	2.81	0.48
3:B:155:LYS:HB3	3:B:219:ARG:HB3	1.96	0.48
3:B:213:VAL:HG21	3:B:226:TYR:HD2	1.79	0.48
3:B:229:LYS:HD2	3:B:237:HIS:HB2	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:23:DT:H1'	1:C:24:DG:H5''	1.96	0.48
3:B:206:ARG:HA	3:B:210:GLN:OE1	2.13	0.48
3:A:25:THR:C	3:A:27:ALA:N	2.70	0.47
3:A:42:PHE:HE1	3:A:115:PHE:HE1	1.62	0.47
3:B:116:ILE:HG13	3:B:133:PRO:HD3	1.96	0.47
3:B:145:ARG:HH21	3:B:167:GLU:CD	2.21	0.47
3:B:185:ILE:HA	3:B:188:ILE:HG12	1.96	0.47
3:A:183:LEU:HD12	3:A:240:PHE:CE1	2.50	0.47
3:A:203:ASP:O	3:A:204:PHE:C	2.56	0.47
3:A:190:GLU:O	3:A:193:GLN:N	2.39	0.47
3:A:132:ASN:ND2	3:A:134:TYR:N	2.62	0.47
3:A:139:PHE:C	3:A:141:GLY:N	2.72	0.47
3:A:32:SER:O	3:A:36:LYS:HG2	2.14	0.47
3:A:68:ALA:O	3:A:73:LEU:N	2.47	0.47
3:A:117:LYS:HB2	3:B:112:PHE:CD2	2.49	0.47
3:A:160:PRO:O	3:A:164:ARG:CG	2.57	0.47
3:A:230:LYS:CB	3:A:235:THR:HA	2.45	0.47
3:B:162:ALA:HA	3:B:212:CYS:HB3	1.97	0.47
3:B:196:GLN:O	3:B:198:TYR:N	2.47	0.47
1:C:17:DA:H2''	1:C:18:DA:H5''	1.96	0.47
3:B:174:LYS:O	3:B:175:PRO:C	2.58	0.47
3:B:142:LEU:O	3:B:145:ARG:HB2	2.15	0.47
1:C:21:DC:C2'	1:C:22:DT:H71	2.38	0.47
3:B:35:GLN:CG	3:B:85:ALA:HB1	2.45	0.46
3:B:213:VAL:HG13	3:B:224:LEU:HD23	1.97	0.46
2:D:63:DC:H2''	2:D:64:DT:C7	2.46	0.46
3:A:29:TYR:HA	3:A:36:LYS:NZ	2.31	0.46
3:A:185:ILE:HG12	3:A:236:THR:O	2.16	0.46
3:A:229:LYS:O	3:A:236:THR:CG2	2.63	0.46
3:A:207:ARG:HA	3:A:207:ARG:HD3	1.57	0.46
3:A:32:SER:OG	3:A:35:GLN:CG	2.60	0.46
3:A:46:ILE:O	3:A:46:ILE:HG22	2.15	0.46
3:A:213:VAL:HG22	3:A:224:LEU:HD23	1.97	0.46
3:A:216:ILE:HA	3:A:220:THR:HG23	1.97	0.46
3:A:236:THR:OG1	3:A:237:HIS:CD2	2.69	0.46
3:A:47:ARG:HG2	3:A:47:ARG:NH1	2.31	0.46
3:B:173:ARG:HB2	3:B:173:ARG:HH11	1.81	0.46
3:A:70:ILE:O	3:A:71:PHE:CD2	2.69	0.46
3:A:191:ARG:O	3:A:192:TYR:HD1	1.98	0.46
3:B:173:ARG:HB2	3:B:173:ARG:NH1	2.30	0.46
3:B:230:LYS:HD3	3:B:233:ARG:O	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:18:DA:H5'	1:C:18:DA:C8	2.51	0.46
2:D:60:DA:C2'	2:D:61:DC:C5'	2.93	0.46
3:A:120:HIS:CE1	3:A:127:TYR:CE1	3.04	0.46
3:A:183:LEU:HD12	3:A:240:PHE:HE1	1.81	0.46
3:B:37:ARG:CZ	3:B:71:PHE:CD2	2.99	0.46
3:A:192:TYR:O	3:A:193:GLN:C	2.59	0.45
3:B:31:LEU:CD1	3:B:39:LEU:HD22	2.45	0.45
3:B:89:PHE:C	3:B:90:ALA:O	2.54	0.45
3:B:120:HIS:CD2	3:B:122:PRO:HD3	2.51	0.45
3:B:146:PHE:O	3:B:150:ARG:HA	2.16	0.45
1:C:24:DG:C2'	1:C:25:DT:H5'	2.31	0.45
3:A:117:LYS:HE3	3:B:112:PHE:HE2	1.81	0.45
3:A:174:LYS:C	3:A:176:ASP:N	2.74	0.45
3:B:172:TYR:CD1	3:B:181:VAL:HG21	2.43	0.45
3:A:122:PRO:HG2	3:B:102:ASP:C	2.40	0.45
3:A:169:LEU:CD1	3:A:224:LEU:HD11	2.46	0.45
3:A:190:GLU:C	3:A:192:TYR:N	2.75	0.45
3:B:184:LYS:HD2	3:B:237:HIS:HD2	1.77	0.45
1:C:5:DT:OP1	3:B:205:ARG:NH2	2.50	0.45
3:A:84:GLN:HE21	3:A:84:GLN:C	2.25	0.45
3:B:33:ARG:O	3:B:37:ARG:HG3	2.17	0.45
3:B:203:ASP:O	3:B:205:ARG:N	2.49	0.45
3:A:182:SER:HB2	3:A:239:VAL:HG22	1.99	0.45
3:B:30:SER:O	3:B:31:LEU:C	2.60	0.45
2:D:41:DG:O6	3:A:206:ARG:NH2	2.45	0.45
3:A:102:ASP:HB2	3:A:106:GLU:O	2.16	0.45
3:B:139:PHE:HA	3:B:142:LEU:HB3	1.99	0.45
3:B:213:VAL:HA	3:B:224:LEU:HD23	1.98	0.45
3:A:84:GLN:HE21	3:A:84:GLN:CA	2.29	0.45
3:A:188:ILE:C	3:A:190:GLU:N	2.73	0.45
3:A:190:GLU:O	3:A:192:TYR:N	2.49	0.45
3:B:78:ALA:O	3:B:81:ASP:N	2.50	0.45
3:B:155:LYS:HE3	3:B:155:LYS:HB2	1.84	0.45
3:B:136:ILE:O	3:B:137:PRO:C	2.59	0.45
3:B:67:TYR:CD2	3:B:68:ALA:N	2.84	0.45
3:A:145:ARG:O	3:A:146:PHE:C	2.58	0.45
3:A:109:TYR:HB2	3:A:110:GLU:H	1.55	0.44
3:A:176:ASP:N	3:A:176:ASP:OD1	2.49	0.44
3:B:133:PRO:HA	3:B:136:ILE:HG13	1.98	0.44
3:B:136:ILE:HB	3:B:137:PRO:CD	2.47	0.44
3:B:189:ILE:HG22	3:B:190:GLU:N	2.31	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:117:LYS:HB2	3:B:112:PHE:CE2	2.52	0.44
3:A:135:LEU:C	3:A:137:PRO:HD2	2.41	0.44
3:B:173:ARG:C	3:B:174:LYS:HG3	2.41	0.44
1:C:22:DT:H2''	1:C:23:DT:C5'	2.37	0.44
3:A:201:MET:O	3:A:204:PHE:HB3	2.18	0.44
3:B:187:TRP:C	3:B:189:ILE:N	2.76	0.44
1:C:28:DC:C5'	1:C:28:DC:H6	2.30	0.44
2:D:62:DA:N3	2:D:63:DC:O4'	2.51	0.44
3:A:41:LEU:O	3:A:44:ASP:N	2.50	0.44
3:A:116:ILE:HG13	3:A:132:ASN:CA	2.45	0.44
3:A:170:CYS:O	3:A:172:TYR:N	2.51	0.44
3:B:25:THR:C	3:B:27:ALA:N	2.74	0.44
3:B:198:TYR:CE1	3:B:207:ARG:CB	2.99	0.44
3:A:165:LEU:HD23	3:A:169:LEU:HG	1.99	0.44
3:B:164:ARG:O	3:B:167:GLU:N	2.50	0.44
3:B:165:LEU:O	3:B:165:LEU:HG	2.18	0.44
3:A:83:ARG:HB2	3:A:127:TYR:OH	2.18	0.44
3:A:99:PRO:HG2	3:A:100:GLU:N	2.23	0.44
3:A:120:HIS:ND1	3:A:127:TYR:CD1	2.85	0.44
3:A:222:MET:HE2	3:A:242:PHE:CG	2.53	0.43
3:B:23:ASP:C	3:B:25:THR:N	2.75	0.43
2:D:35:DA:H1'	2:D:36:DA:C5'	2.46	0.43
3:A:97:TYR:HE2	3:A:107:LYS:NZ	2.17	0.43
2:D:58:DT:H1'	2:D:59:DC:H5''	1.99	0.43
3:A:33:ARG:O	3:A:37:ARG:HG3	2.17	0.43
3:A:45:GLN:C	3:A:47:ARG:N	2.76	0.43
3:A:169:LEU:HB3	3:A:242:PHE:CE2	2.53	0.43
3:A:200:ARG:HD3	3:A:203:ASP:OD2	2.18	0.43
3:B:112:PHE:CE2	3:B:134:TYR:CD2	3.06	0.43
2:D:54:DA:H2''	2:D:55:DT:H5''	1.98	0.43
2:D:55:DT:C2'	2:D:56:DT:H5''	2.47	0.43
3:B:83:ARG:HG3	3:B:84:GLN:N	2.33	0.43
2:D:62:DA:H1'	2:D:63:DC:C5'	2.49	0.43
3:A:185:ILE:HD11	3:A:238:ILE:HG12	2.00	0.43
3:B:146:PHE:CE1	3:B:151:LEU:HD23	2.54	0.43
3:B:195:PRO:HB2	3:B:198:TYR:CD2	2.54	0.43
1:C:21:DC:C2'	1:C:22:DT:H5''	2.36	0.43
2:D:57:DG:H2''	2:D:58:DT:C5'	2.48	0.43
3:A:109:TYR:N	3:A:109:TYR:CD1	2.87	0.43
3:A:211:VAL:HG12	3:A:212:CYS:N	2.32	0.43
3:B:24:LEU:HD13	3:B:138:PHE:HB2	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:160:PRO:C	3:B:163:MET:HG2	2.43	0.43
1:C:7:DT:H2''	1:C:8:DG:C8	2.54	0.43
3:A:139:PHE:O	3:A:142:LEU:N	2.41	0.43
3:A:180:ILE:HG12	3:A:241:SER:OG	2.18	0.43
3:B:100:GLU:CG	3:B:101:GLU:N	2.77	0.43
3:A:201:MET:HB3	3:A:202:PRO:HD3	2.01	0.42
3:B:40:TYR:C	3:B:42:PHE:N	2.76	0.42
3:B:121:SER:O	3:B:123:SER:N	2.52	0.42
3:B:146:PHE:CD2	3:B:152:SER:HB2	2.51	0.42
3:B:32:SER:HA	3:B:156:GLU:CD	2.44	0.42
3:B:121:SER:O	3:B:122:PRO:C	2.61	0.42
3:B:164:ARG:HB3	3:B:192:TYR:CD1	2.53	0.42
3:B:220:THR:HB	3:B:221:PRO:HD2	2.00	0.42
3:B:169:LEU:HD21	3:B:240:PHE:CD1	2.54	0.42
3:A:38:MET:O	3:A:39:LEU:C	2.61	0.42
3:B:112:PHE:N	3:B:112:PHE:CD1	2.87	0.42
1:C:6:DG:O6	2:D:61:DC:N4	2.42	0.42
2:D:44:DA:H1'	2:D:45:DA:H5''	2.01	0.42
2:D:45:DA:C2'	2:D:46:DG:C5'	2.97	0.42
3:A:73:LEU:CD2	3:A:77:GLU:OE1	2.63	0.42
1:C:30:DC:C6	1:C:31:DT:H72	2.55	0.42
3:A:159:ASN:OD1	3:A:162:ALA:N	2.50	0.42
3:B:24:LEU:CD1	3:B:138:PHE:HB2	2.49	0.42
3:B:33:ARG:HH21	3:B:34:ASP:CG	2.27	0.42
3:B:226:TYR:CD1	3:B:226:TYR:C	2.97	0.42
2:D:48:DT:H2''	2:D:49:DT:C6	2.55	0.42
3:A:79:SER:O	3:A:80:LYS:C	2.63	0.42
3:A:172:TYR:CD1	3:A:181:VAL:HG21	2.54	0.42
3:A:209:LEU:HD13	3:A:226:TYR:CE1	2.55	0.42
3:B:93:GLU:HG2	3:B:113:PRO:HA	2.01	0.42
3:B:133:PRO:HA	3:B:136:ILE:CG1	2.49	0.42
3:A:168:SER:CB	3:A:192:TYR:OH	2.68	0.42
3:B:30:SER:OG	3:B:31:LEU:N	2.52	0.42
3:B:178:SER:HA	3:B:243:ARG:HA	2.01	0.42
3:B:165:LEU:HD11	3:B:240:PHE:CE1	2.55	0.42
3:B:174:LYS:NZ	3:B:174:LYS:HB3	2.35	0.42
3:A:159:ASN:O	3:A:161:TYR:N	2.53	0.41
3:A:170:CYS:O	3:A:171:GLN:C	2.62	0.41
3:A:223:ARG:HA	3:A:223:ARG:HD2	1.78	0.41
2:D:39:DG:O6	3:A:200:ARG:NH2	2.54	0.41
3:A:99:PRO:CG	3:A:100:GLU:H	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:223:ARG:HD3	3:A:245:ILE:HB	2.01	0.41
3:B:32:SER:OG	3:B:35:GLN:HG3	2.20	0.41
3:B:117:LYS:HD3	3:B:130:HIS:CE1	2.54	0.41
3:B:129:VAL:C	3:B:130:HIS:HD2	2.28	0.41
1:C:11:DA:C6	1:C:12:DA:C6	3.08	0.41
3:B:206:ARG:HE	3:B:206:ARG:HB3	1.70	0.41
3:A:100:GLU:O	3:A:102:ASP:N	2.54	0.41
3:A:198:TYR:N	3:A:198:TYR:CD2	2.89	0.41
3:B:187:TRP:O	3:B:191:ARG:HB2	2.20	0.41
1:C:1:DT:C2'	1:C:2:DT:H5'	2.36	0.41
3:A:241:SER:O	3:A:243:ARG:NH2	2.53	0.41
3:B:65:ALA:O	3:B:66:LYS:C	2.62	0.41
1:C:12:DA:H2''	1:C:13:DT:H5'	2.02	0.41
3:A:78:ALA:O	3:A:82:ILE:HG12	2.21	0.41
3:A:170:CYS:O	3:A:173:ARG:N	2.50	0.41
3:B:25:THR:CG2	3:B:145:ARG:NE	2.84	0.41
3:B:23:ASP:O	3:B:24:LEU:C	2.64	0.41
3:A:93:GLU:HA	3:A:113:PRO:HA	2.03	0.41
3:A:188:ILE:O	3:A:190:GLU:N	2.54	0.41
3:A:202:PRO:O	3:A:205:ARG:HB2	2.20	0.41
3:B:164:ARG:C	3:B:166:TYR:H	2.29	0.41
1:C:11:DA:H2''	1:C:12:DA:OP2	2.21	0.41
1:C:16:DA:H2	2:D:52:DA:C2	2.40	0.41
3:A:159:ASN:HA	3:A:160:PRO:HD2	1.80	0.41
3:A:169:LEU:HD11	3:A:224:LEU:HD21	2.03	0.41
3:A:242:PHE:CD1	3:A:242:PHE:C	2.98	0.41
3:B:115:PHE:CD2	3:B:129:VAL:HG13	2.55	0.41
3:A:43:VAL:O	3:A:47:ARG:HG3	2.21	0.40
2:D:45:DA:H1'	2:D:46:DG:C5'	2.49	0.40
2:D:51:DT:H1'	2:D:52:DA:H5'	2.03	0.40
2:D:55:DT:H1'	2:D:56:DT:H5''	2.02	0.40
2:D:63:DC:H2''	2:D:64:DT:OP2	2.21	0.40
3:A:31:LEU:HB2	3:A:36:LYS:HD2	2.02	0.40
3:A:185:ILE:HG21	3:A:201:MET:HE1	2.03	0.40
3:B:99:PRO:O	3:B:102:ASP:OD1	2.38	0.40
3:A:33:ARG:O	3:A:34:ASP:C	2.65	0.40
3:A:145:ARG:HA	3:A:149:PHE:HD1	1.86	0.40
3:B:173:ARG:CZ	3:B:173:ARG:HB3	2.51	0.40
2:D:45:DA:C1'	2:D:46:DG:H5''	2.50	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	
3	A	218/266 (82%)	158 (72%)	40 (18%)	20 (9%)	0	2
3	B	214/266 (80%)	158 (74%)	35 (16%)	21 (10%)	0	2
All	All	432/532 (81%)	316 (73%)	75 (17%)	41 (10%)	0	2

All (41) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	A	91	GLY
3	A	99	PRO
3	B	68	ALA
3	B	147	THR
3	B	209	LEU
3	B	211	VAL
3	A	46	ILE
3	A	101	GLU
3	A	191	ARG
3	A	209	LEU
3	B	31	LEU
3	B	125	GLY
3	B	165	LEU
3	B	173	ARG
3	A	39	LEU
3	A	72	GLY
3	A	82	ILE
3	A	171	GLN
3	A	175	PRO
3	B	22	ASN
3	B	67	TYR
3	B	204	PHE
3	B	210	GLN
3	B	215	GLU
3	A	79	SER

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Mol	Chain	Res	Type
3	A	81	ASP
3	A	201	MET
3	B	45	GLN
3	B	64	VAL
3	B	197	SER
3	B	202	PRO
3	B	246	THR
3	A	140	ILE
3	A	152	SER
3	A	193	GLN
3	A	211	VAL
3	B	23	ASP
3	B	122	PRO
3	A	125	GLY
3	B	188	ILE
3	A	58	GLY

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	
3	A	188/238 (79%)	165 (88%)	23 (12%)	5	18
3	B	189/238 (79%)	174 (92%)	15 (8%)	11	34
All	All	377/476 (79%)	339 (90%)	38 (10%)	7	25

All (38) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	A	36	LYS
3	A	39	LEU
3	A	42	PHE
3	A	69	GLU
3	A	84	GLN
3	A	89	PHE
3	A	100	GLU

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Mol	Chain	Res	Type
3	A	117	LYS
3	A	126	LEU
3	A	132	ASN
3	A	135	LEU
3	A	145	ARG
3	A	152	SER
3	A	165	LEU
3	A	176	ASP
3	A	189	ILE
3	A	193	GLN
3	A	207	ARG
3	A	218	SER
3	A	220	THR
3	A	245	ILE
3	A	246	THR
3	A	247	SER
3	B	80	LYS
3	B	82	ILE
3	B	89	PHE
3	B	98	ARG
3	B	100	GLU
3	B	126	LEU
3	B	135	LEU
3	B	154	THR
3	B	164	ARG
3	B	165	LEU
3	B	185	ILE
3	B	191	ARG
3	B	215	GLU
3	B	225	SER
3	B	226	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
3	A	35	GLN
3	A	45	GLN
3	A	84	GLN
3	A	132	ASN
3	A	148	GLN
3	A	171	GLN
3	A	196	GLN

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Mol	Chain	Res	Type
3	A	237	HIS
3	B	84	GLN
3	B	130	HIS
3	B	148	GLN
3	B	159	ASN

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 oligosaccharides 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, 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	C	33/33 (100%)	-1.67	0 100 100	22, 65, 95, 99	0
2	D	33/33 (100%)	-1.65	0 100 100	39, 59, 103, 109	0
3	A	222/266 (83%)	-1.55	0 100 100	3, 33, 66, 79	0
3	B	218/266 (81%)	-1.53	0 100 100	5, 41, 69, 80	0
All	All	506/598 (84%)	-1.56	0 100 100	3, 40, 75, 109	0

There are no RSRZ outliers to report.

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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.