



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 01:58 AM UTC

PDB ID : 4PXI / pdb_00004pxi
Title : Elucidation of the Structural and Functional Mechanism of Action of the TetR Family Protein, CprB from *S. coelicolor* A3(2)
Authors : Hussain, B.; Ruchika, B.; Aruna, B.; Ruchi, A.
Deposited on : 2014-03-24
Resolution : 3.20 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:

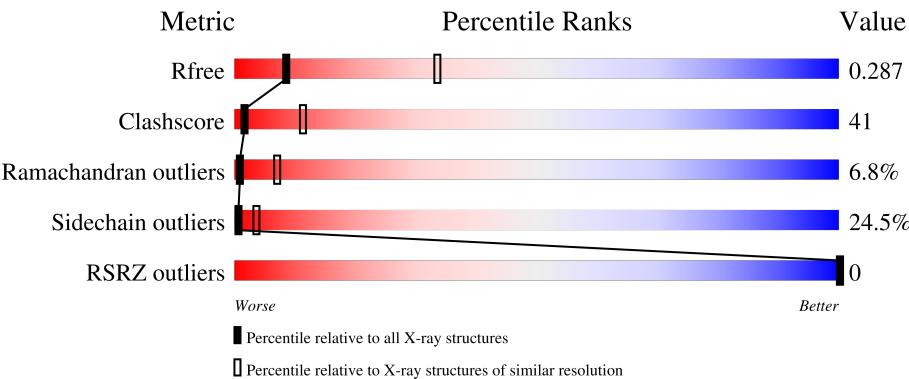
MolProbity : 4-5-2 with Phenix2.0
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
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 3.20 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	1466 (3.20-3.20)
Clashscore	190562	1573 (3.20-3.20)
Ramachandran outliers	187476	1548 (3.20-3.20)
Sidechain outliers	187428	1547 (3.20-3.20)
RSRZ outliers	180081	1466 (3.20-3.20)

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	215	34%37%15%•10%
1	B	215	35%50%8%•5%
1	C	215	32%44%16%•7%
1	D	215	34%43%12%•10%
2	E	22	14%41%36%9%

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Mol	Chain	Length	Quality of chain
3	F	22	 A horizontal bar chart showing the quality of chain F. The bar is divided into three segments: a green segment on the left labeled '5%', a yellow segment in the middle labeled '41%', and an orange segment on the right labeled '55%'. The segments are separated by thin black lines.

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6886 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 CprB.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	194	Total	C	N	O	S	0	0	0
			1480	930	276	269	5			
1	B	204	Total	C	N	O	S	0	0	0
			1518	950	277	285	6			
1	C	200	Total	C	N	O	S	0	0	0
			1535	961	295	274	5			
1	D	193	Total	C	N	O	S	0	0	0
			1497	939	279	273	6			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	20	Total	C	N	O	P	0	0	0
			405	194	73	119	19			

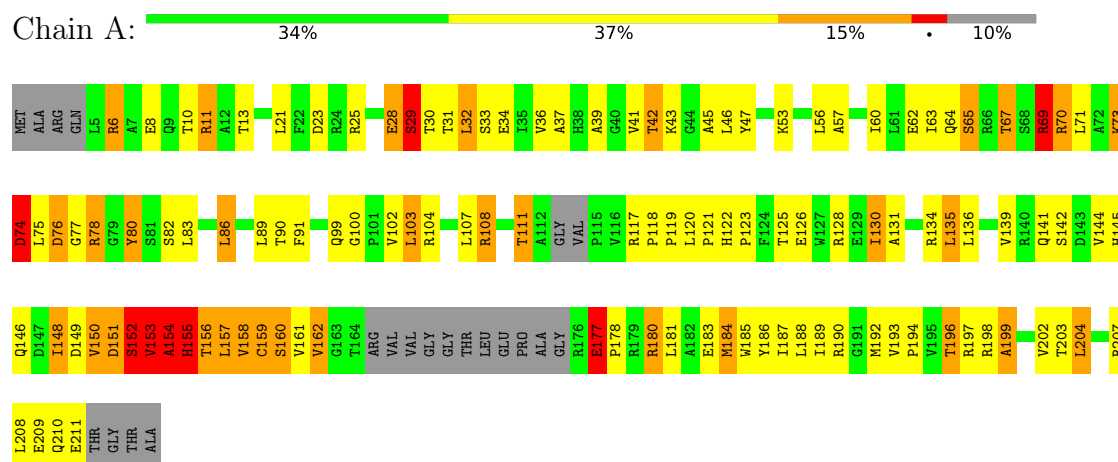
- Molecule 3 is a DNA chain called DNA (5'-D(*AP*TP*AP*AP*AP*CP*GP*GP*GP*GP*CP*GP*TP*CP*CP*CP*GP*TP*AP*TP*GP*T)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	F	22	Total	C	N	O	P	0	0	0
			451	215	85	130	21			

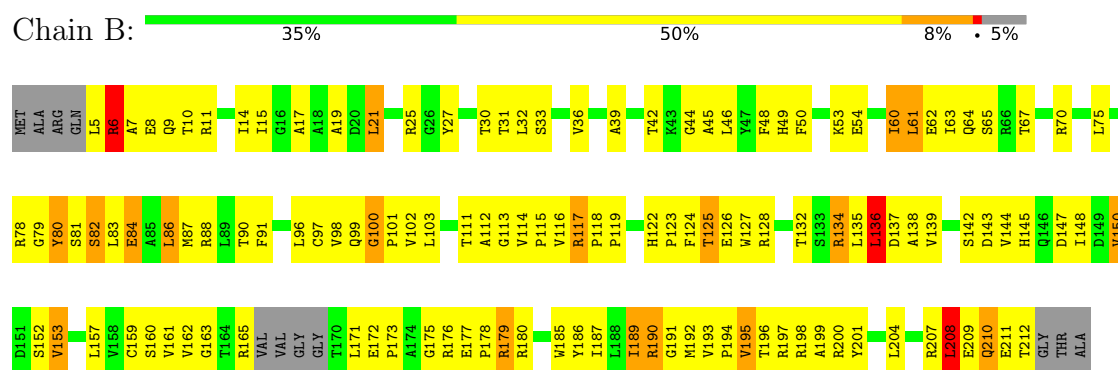
3 Residue-property plots

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

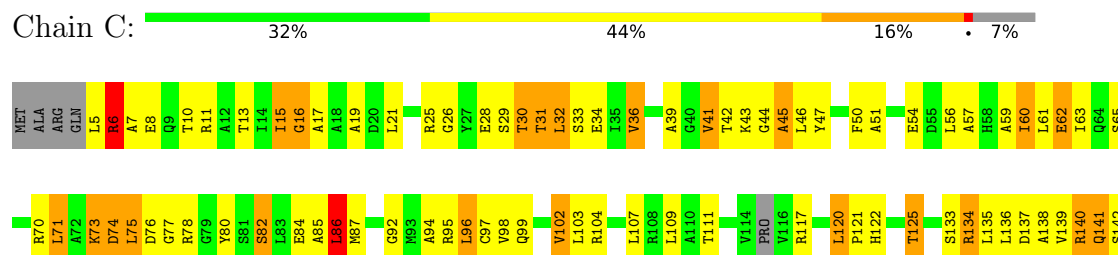
• Molecule 1: CprB

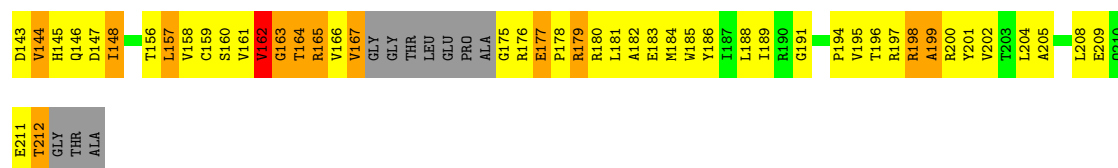


• Molecule 1: CprB

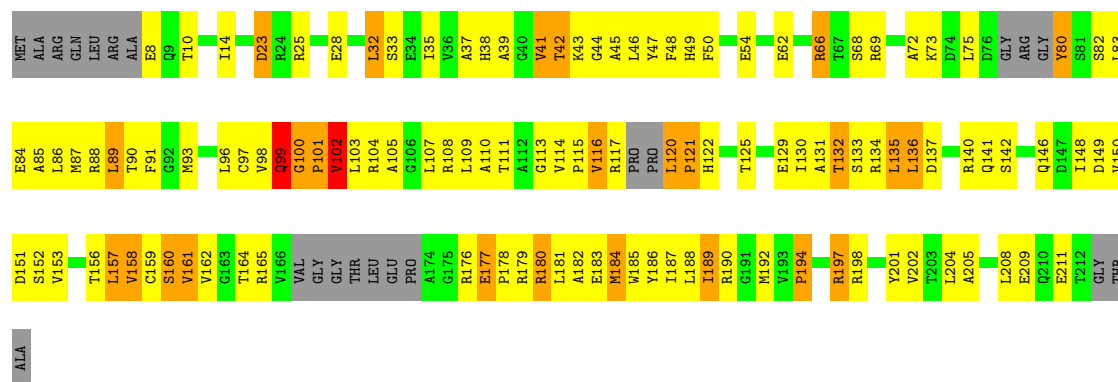


• Molecule 1: CprB

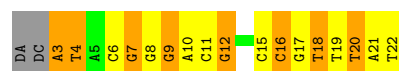




• Molecule 1: CprB



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



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



4 Data and refinement statistics

Property	Value	Source
Space group	P 32	Depositor
Cell constants a, b, c, α , β , γ	149.06Å 149.06Å 69.07Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	60.90 – 3.20 60.90 – 3.20	Depositor EDS
% Data completeness (in resolution range)	99.6 (60.90-3.20) 99.6 (60.90-3.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.13	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.55 (at 3.19Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.220 , 0.294 0.218 , 0.287	Depositor DCC
R_{free} test set	1163 reflections (4.10%)	wwPDB-VP
Wilson B-factor (Å ²)	74.4	Xtriage
Anisotropy	0.409	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 98.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.034 for -h,-k,l 0.428 for h,-h-k,-l 0.035 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6886	wwPDB-VP
Average B, all atoms (Å ²)	88.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.41% 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 i

5.1 Standard geometry i

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	0.95	2/1505 (0.1%)	1.19	9/2043 (0.4%)
1	B	0.92	0/1545	1.25	12/2104 (0.6%)
1	C	0.89	0/1559	1.20	6/2111 (0.3%)
1	D	0.87	0/1519	1.08	3/2056 (0.1%)
2	E	0.57	0/453	1.46	11/697 (1.6%)
3	F	0.54	0/506	1.55	16/780 (2.1%)
All	All	0.87	2/7087 (0.0%)	1.24	57/9791 (0.6%)

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	A	0	1
1	D	0	1
All	All	0	2

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	155	HIS	CG-CD2	5.88	1.42	1.35
1	A	118	PRO	CA-C	5.73	1.55	1.51

All (57) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	3	DA	P-O3'-C3'	12.24	138.56	120.20
3	F	1	DA	O3'-P-O5'	10.81	120.22	104.00
3	F	3	DA	P-O3'-C3'	10.62	136.13	120.20
1	C	6	ARG	N-CA-C	-10.13	100.33	112.89
3	F	1	DA	C2'-C3'-O3'	-9.59	97.12	111.50
1	C	165	ARG	N-CA-C	-9.07	102.22	113.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	F	21	DG	P-O3'-C3'	8.86	133.49	120.20
1	B	172	GLU	CA-C-N	7.95	129.78	119.84
1	B	172	GLU	C-N-CA	7.95	129.78	119.84
3	F	13	DT	C2'-C3'-O3'	-7.83	99.76	111.50
3	F	10	DG	P-O3'-C3'	-7.35	109.17	120.20
3	F	3	DA	O3'-P-O5'	7.33	114.99	104.00
2	E	4	DT	P-O3'-C3'	7.28	131.12	120.20
2	E	20	DT	P-O3'-C3'	7.27	131.10	120.20
1	A	162	VAL	N-CA-C	7.19	118.82	110.62
2	E	9	DG	O3'-P-O5'	-7.17	93.25	104.00
1	A	57	ALA	N-CA-C	-6.76	103.60	110.97
3	F	18	DT	P-O3'-C3'	6.60	130.10	120.20
3	F	8	DG	O3'-P-O5'	-6.51	94.24	104.00
2	E	15	DC	C2'-C3'-O3'	-6.50	101.75	111.50
2	E	12	DG	P-O3'-C3'	-6.46	110.52	120.20
3	F	7	DG	P-O3'-C3'	-6.40	110.60	120.20
1	C	125	THR	N-CA-C	-6.39	104.39	111.36
3	F	13	DT	P-O3'-C3'	6.36	129.74	120.20
1	D	100	GLY	CA-C-N	6.33	127.75	119.84
1	D	100	GLY	C-N-CA	6.33	127.75	119.84
2	E	16	DC	O3'-P-O5'	-6.25	94.63	104.00
1	B	100	GLY	CA-C-N	6.24	125.93	119.56
1	B	100	GLY	C-N-CA	6.24	125.93	119.56
3	F	16	DC	O3'-P-O5'	-6.22	94.67	104.00
3	F	5	DA	C2'-C3'-O3'	-5.98	102.53	111.50
1	C	63	ILE	CB-CA-C	-5.98	104.32	111.97
1	A	80	TYR	N-CA-C	5.95	118.89	110.50
1	D	152	SER	N-CA-C	-5.94	104.81	111.28
3	F	20	DT	C2'-C3'-O3'	5.91	120.36	111.50
1	A	153	VAL	N-CA-C	5.89	121.60	109.34
3	F	19	DA	P-O3'-C3'	5.77	128.85	120.20
1	B	70	ARG	N-CA-C	-5.75	106.22	113.18
2	E	18	DT	O3'-P-O5'	-5.68	95.49	104.00
3	F	14	DC	O3'-P-O5'	-5.56	95.66	104.00
1	B	125	THR	N-CA-C	-5.52	106.25	112.87
1	A	154	ALA	N-CA-C	5.39	122.27	110.80
1	A	154	ALA	CA-C-N	5.33	127.36	120.44
1	A	154	ALA	C-N-CA	5.33	127.36	120.44
2	E	9	DG	P-O3'-C3'	-5.25	112.32	120.20
1	B	114	VAL	CA-C-N	5.22	125.96	120.11
1	B	114	VAL	C-N-CA	5.22	125.96	120.11
2	E	3	DA	N9-C1'-C2'	5.19	121.29	113.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	150	VAL	N-CA-C	5.17	115.61	110.23
1	A	117	ARG	CA-C-N	5.14	123.41	119.66
1	A	117	ARG	C-N-CA	5.14	123.41	119.66
1	C	61	LEU	N-CA-C	-5.13	105.38	111.69
1	B	189	ILE	N-CA-C	-5.13	102.98	109.80
1	C	45	ALA	N-CA-C	-5.09	105.81	111.36
1	B	48	PHE	N-CA-C	-5.05	105.67	111.07
2	E	7	DG	C4'-C3'-O3'	-5.02	102.47	110.00
1	B	45	ALA	N-CA-C	-5.01	105.53	111.69

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	152	SER	Peptide
1	D	125	THR	Peptide

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	1480	0	1469	124	0
1	B	1518	0	1464	117	0
1	C	1535	0	1542	124	0
1	D	1497	0	1509	120	0
2	E	405	0	227	52	0
3	F	451	0	249	65	0
All	All	6886	0	6460	544	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 41.

All (544) 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:C:144:VAL:HG12	1:C:145:HIS:H	1.16	1.09

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:177:GLU:H	1:A:178:PRO:HD2	1.20	1.04
1:A:154:ALA:HA	1:A:157:LEU:HD13	1.40	1.03
1:C:163:GLY:HA3	1:C:166:VAL:HB	1.06	1.02
1:A:151:ASP:HB3	1:A:153:VAL:HG22	1.41	1.02
1:D:42:THR:HG23	1:D:45:ALA:HB2	1.41	1.02
1:B:86:LEU:HB2	1:B:134:ARG:HG2	1.41	1.02
1:D:180:ARG:HG3	1:D:180:ARG:HH11	1.19	1.01
1:C:163:GLY:HA3	1:C:166:VAL:CB	1.91	1.00
1:C:163:GLY:CA	1:C:166:VAL:HB	1.90	0.99
2:E:10:DA:H2''	2:E:11:DC:H5''	1.43	0.99
1:B:80:TYR:HB3	1:B:84:GLU:HB2	1.42	0.99
1:C:33:SER:HA	1:C:36:VAL:CG2	1.94	0.97
1:D:45:ALA:CA	3:F:1:DA:H2'	1.93	0.97
1:D:45:ALA:HA	3:F:1:DA:H2'	1.46	0.95
1:C:142:SER:HB3	1:C:197:ARG:HH21	1.29	0.95
1:C:175:GLY:HA2	1:C:176:ARG:HB3	1.48	0.93
1:C:33:SER:HA	1:C:36:VAL:HG23	1.49	0.92
3:F:9:DG:H4'	3:F:9:DG:OP1	1.66	0.92
1:C:26:GLY:O	1:C:30:THR:HG23	1.70	0.91
1:B:210:GLN:CD	1:B:210:GLN:H	1.78	0.91
1:B:189:ILE:O	1:B:190:ARG:HB2	1.71	0.91
1:C:144:VAL:HG12	1:C:145:HIS:N	1.84	0.89
2:E:21:DA:H2''	2:E:22:DT:O5'	1.72	0.89
3:F:19:DA:H1'	3:F:20:DT:H5'	1.55	0.88
1:A:42:THR:HG23	1:A:45:ALA:HB2	1.56	0.88
1:B:7:ALA:HB2	3:F:4:DA:H5''	1.55	0.87
2:E:17:DG:H1	3:F:6:DC:H42	1.22	0.87
1:A:196:THR:HG22	1:A:197:ARG:N	1.89	0.86
1:A:86:LEU:HD12	1:A:134:ARG:HB3	1.57	0.86
1:A:154:ALA:HA	1:A:157:LEU:CD1	2.06	0.86
3:F:7:DG:H2''	3:F:8:DG:C8	2.13	0.84
3:F:2:DT:H2''	3:F:3:DA:H5'	1.58	0.83
1:B:86:LEU:HD23	1:B:134:ARG:HB3	1.59	0.83
2:E:16:DC:H2''	2:E:17:DG:C8	2.14	0.83
1:B:190:ARG:HE	1:B:198:ARG:HH21	1.23	0.83
1:B:132:THR:O	1:B:136:LEU:HB3	1.79	0.82
1:C:134:ARG:HG3	1:C:134:ARG:HH11	1.42	0.82
1:D:42:THR:HG23	1:D:45:ALA:CB	2.09	0.82
2:E:11:DC:H2''	2:E:12:DG:C8	2.15	0.81
1:B:207:ARG:O	1:B:208:LEU:HB2	1.79	0.81
1:A:180:ARG:HB2	1:A:180:ARG:HH11	1.47	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:69:ARG:NH1	1:D:69:ARG:HB3	1.97	0.80
1:B:159:CYS:O	1:B:162:VAL:HG22	1.80	0.79
1:C:111:THR:HG21	1:D:111:THR:HG21	1.63	0.79
2:E:21:DA:H2''	2:E:22:DT:H3'	1.64	0.79
1:A:42:THR:HG21	2:E:4:DT:OP1	1.82	0.78
1:B:139:VAL:HG22	1:B:144:VAL:HG21	1.66	0.78
1:B:163:GLY:HA3	1:B:180:ARG:NH2	2.00	0.77
2:E:7:DG:N2	3:F:16:DC:N3	2.29	0.77
3:F:6:DC:H2''	3:F:7:DG:C8	2.20	0.77
1:C:96:LEU:HB2	1:C:103:LEU:HD12	1.66	0.76
1:B:96:LEU:HB2	1:B:103:LEU:HD12	1.67	0.76
2:E:6:DC:H2''	2:E:7:DG:C8	2.20	0.76
2:E:7:DG:H1	3:F:16:DC:H42	1.32	0.76
1:C:145:HIS:CD2	1:D:190:ARG:HH12	2.03	0.76
2:E:6:DC:C2'	2:E:7:DG:C8	2.69	0.76
1:A:42:THR:HG23	1:A:45:ALA:CB	2.17	0.75
1:B:204:LEU:O	1:B:208:LEU:HB2	1.86	0.75
1:C:96:LEU:HB2	1:C:103:LEU:CD1	2.16	0.75
1:C:159:CYS:HA	1:C:162:VAL:HB	1.67	0.75
2:E:17:DG:N2	3:F:6:DC:N3	2.33	0.74
3:F:2:DT:H6	3:F:2:DT:H5''	1.53	0.74
1:B:157:LEU:O	1:B:161:VAL:HG23	1.88	0.73
1:D:87:MET:HE1	1:D:205:ALA:HB2	1.71	0.73
1:B:80:TYR:CB	1:B:84:GLU:HB2	2.18	0.73
1:A:145:HIS:CE1	1:A:194:PRO:HA	2.22	0.73
1:A:153:VAL:O	1:A:155:HIS:HD2	1.72	0.73
1:D:80:TYR:OH	1:D:88:ARG:NH1	2.22	0.73
1:D:86:LEU:HD21	1:D:135:LEU:HD13	1.70	0.73
1:D:69:ARG:HB3	1:D:69:ARG:HH11	1.53	0.73
2:E:17:DG:H1	3:F:6:DC:N4	1.86	0.73
1:A:183:GLU:HA	1:A:186:TYR:HD2	1.53	0.73
1:C:142:SER:HB3	1:C:197:ARG:NH2	2.04	0.73
1:A:25:ARG:HB2	1:A:30:THR:OG1	1.89	0.73
1:D:32:LEU:HA	1:D:35:ILE:HD12	1.71	0.73
2:E:10:DA:C2'	2:E:11:DC:H5''	2.18	0.72
1:C:94:ALA:O	1:C:98:VAL:HG23	1.89	0.72
1:A:86:LEU:HD13	1:A:135:LEU:HD13	1.71	0.72
1:A:177:GLU:H	1:A:178:PRO:CD	2.00	0.71
1:A:189:ILE:HG23	1:A:193:VAL:HG21	1.71	0.71
2:E:9:DG:H2''	2:E:10:DA:C8	2.24	0.71
1:C:47:TYR:CD1	3:F:13:DT:H72	2.25	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:86:LEU:HD13	1:A:135:LEU:CD1	2.21	0.71
1:C:76:ASP:C	1:C:78:ARG:H	1.99	0.71
1:C:5:LEU:HD13	3:F:21:DG:OP1	1.90	0.71
1:C:117:ARG:O	1:C:120:LEU:HB3	1.91	0.70
1:D:42:THR:HG21	3:F:2:DT:OP2	1.91	0.70
1:B:139:VAL:HG22	1:B:144:VAL:CG2	2.21	0.70
1:C:181:LEU:HD12	1:C:184:MET:HE3	1.74	0.70
1:A:86:LEU:CD1	1:A:134:ARG:HB3	2.21	0.70
1:B:165:ARG:HA	1:B:171:LEU:HB2	1.73	0.69
1:D:84:GLU:HG3	1:D:204:LEU:HD11	1.74	0.69
3:F:20:DT:H2''	3:F:21:DG:C5	2.28	0.69
1:C:144:VAL:CG1	1:C:145:HIS:H	1.95	0.69
1:A:209:GLU:C	1:A:211:GLU:H	1.99	0.69
1:C:36:VAL:HG13	1:C:46:LEU:HD12	1.75	0.68
1:B:195:VAL:HA	1:B:198:ARG:HD3	1.74	0.68
1:D:82:SER:H	1:D:141:GLN:HE22	1.40	0.68
1:A:155:HIS:CD2	1:A:155:HIS:H	2.09	0.68
1:D:188:LEU:O	1:D:192:MET:HB2	1.92	0.68
1:C:21:LEU:HD22	1:C:25:ARG:NH1	2.09	0.68
1:A:187:ILE:HG23	1:B:191:GLY:HA3	1.75	0.67
1:D:43:LYS:HD2	2:E:18:DT:H72	1.76	0.67
2:E:18:DT:H2'	2:E:19:DT:C7	2.25	0.67
2:E:17:DG:H2''	2:E:18:DT:H5''	1.76	0.67
1:B:145:HIS:O	1:B:148:ILE:HG13	1.95	0.67
1:C:17:ALA:HB3	1:C:39:ALA:HB2	1.75	0.67
1:A:209:GLU:O	1:A:211:GLU:N	2.24	0.67
1:B:86:LEU:O	1:B:90:THR:HG23	1.96	0.66
1:C:145:HIS:HD2	1:D:190:ARG:HH12	1.43	0.66
1:C:175:GLY:HA2	1:C:176:ARG:CB	2.25	0.66
1:A:196:THR:CG2	1:A:197:ARG:N	2.59	0.66
2:E:6:DC:H2''	2:E:7:DG:H8	1.58	0.65
2:E:21:DA:C2'	2:E:22:DT:H3'	2.26	0.65
1:B:211:GLU:O	1:B:211:GLU:HG2	1.94	0.65
1:D:83:LEU:HD13	1:D:83:LEU:O	1.96	0.65
2:E:18:DT:H2'	2:E:19:DT:H72	1.79	0.65
1:A:86:LEU:CD1	1:A:135:LEU:HD13	2.26	0.65
1:D:180:ARG:HG3	1:D:180:ARG:NH1	1.98	0.65
3:F:7:DG:H2''	3:F:8:DG:H8	1.60	0.65
1:C:21:LEU:HD22	1:C:25:ARG:HH11	1.62	0.65
1:C:92:GLY:O	1:C:95:ARG:HB3	1.96	0.65
1:B:210:GLN:CD	1:B:210:GLN:N	2.51	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:186:TYR:CE1	1:D:202:VAL:HG13	2.32	0.65
1:B:19:ALA:HB2	1:B:60:ILE:HD11	1.79	0.64
1:B:189:ILE:O	1:B:190:ARG:CB	2.43	0.64
3:F:9:DG:H2''	3:F:10:DG:C8	2.32	0.64
1:A:151:ASP:OD2	1:A:151:ASP:N	2.29	0.64
1:B:82:SER:HB3	1:B:138:ALA:HB2	1.78	0.64
1:B:21:LEU:HD11	1:B:25:ARG:NH2	2.12	0.64
1:A:177:GLU:N	1:A:178:PRO:HD2	2.04	0.64
1:A:183:GLU:O	1:A:187:ILE:HD12	1.97	0.64
1:C:7:ALA:HB2	2:E:6:DC:H5''	1.79	0.64
2:E:11:DC:H2''	2:E:12:DG:N7	2.13	0.64
1:B:6:ARG:HD3	1:B:6:ARG:N	2.13	0.63
1:D:42:THR:CG2	1:D:45:ALA:HB2	2.24	0.63
1:D:197:ARG:HG3	1:D:197:ARG:HH11	1.62	0.63
2:E:18:DT:H2'	2:E:19:DT:C5	2.34	0.63
1:D:25:ARG:HH21	1:D:25:ARG:HG3	1.61	0.63
1:D:101:PRO:HA	1:D:104:ARG:HH21	1.63	0.63
1:A:144:VAL:HG12	1:A:148:ILE:HD11	1.80	0.63
1:D:180:ARG:HH11	1:D:180:ARG:CG	2.05	0.63
1:B:14:ILE:HG21	1:B:50:PHE:HE2	1.63	0.63
1:B:79:GLY:O	1:B:80:TYR:HB2	1.99	0.62
1:C:19:ALA:HB2	1:C:60:ILE:HD11	1.80	0.62
1:D:86:LEU:CD2	1:D:135:LEU:HD13	2.29	0.62
1:D:88:ARG:HG2	1:D:208:LEU:HD11	1.81	0.62
1:B:64:GLN:O	1:B:67:THR:HG22	1.98	0.62
1:D:113:GLY:O	1:D:115:PRO:HD3	1.98	0.62
1:C:179:ARG:CZ	1:C:179:ARG:HB3	2.28	0.62
1:A:151:ASP:CA	1:A:152:SER:HB3	2.28	0.62
1:A:209:GLU:C	1:A:211:GLU:N	2.57	0.62
1:B:189:ILE:HA	1:B:193:VAL:HG23	1.81	0.62
1:A:150:VAL:O	1:A:152:SER:HB3	2.00	0.62
1:A:183:GLU:HA	1:A:186:TYR:CD2	2.33	0.62
1:D:185:TRP:O	1:D:189:ILE:HG12	2.01	0.61
1:A:125:THR:HG22	1:A:128:ARG:HH21	1.66	0.61
1:D:83:LEU:HD13	1:D:87:MET:HG3	1.81	0.61
1:A:23:ASP:HA	1:A:108:ARG:HE	1.64	0.61
1:B:42:THR:HG21	3:F:6:DC:H3'	1.83	0.61
1:C:134:ARG:HG3	1:C:134:ARG:NH1	2.14	0.61
1:C:159:CYS:O	1:C:160:SER:C	2.43	0.61
1:A:151:ASP:HA	1:A:152:SER:HB3	1.81	0.61
1:D:44:GLY:O	3:F:1:DA:H8	1.84	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:189:ILE:HG22	1:D:198:ARG:HG3	1.83	0.61
1:A:69:ARG:HG2	1:A:69:ARG:HH11	1.66	0.60
1:B:84:GLU:O	1:B:88:ARG:HG2	2.01	0.60
3:F:9:DG:OP1	3:F:9:DG:C4'	2.44	0.60
1:A:157:LEU:HD12	1:A:157:LEU:H	1.66	0.60
1:A:196:THR:HG22	1:A:197:ARG:HD2	1.84	0.60
1:C:145:HIS:HB3	1:C:147:ASP:OD1	2.02	0.60
1:C:42:THR:OG1	2:E:8:DG:OP2	2.18	0.60
1:A:31:THR:HG23	1:A:34:GLU:HG2	1.83	0.60
1:A:107:LEU:HD11	1:A:162:VAL:HG22	1.84	0.60
1:C:144:VAL:CG1	1:C:145:HIS:N	2.56	0.60
1:C:59:ALA:HA	1:C:62:GLU:HB2	1.84	0.60
1:D:120:LEU:HG	1:D:121:PRO:HD2	1.83	0.60
3:F:2:DT:H5''	3:F:2:DT:C6	2.34	0.59
1:B:62:GLU:C	1:B:64:GLN:H	2.10	0.59
2:E:18:DT:H2'	2:E:19:DT:C6	2.38	0.59
1:A:42:THR:CG2	1:A:45:ALA:HB2	2.29	0.59
1:C:31:THR:HG23	1:C:34:GLU:CB	2.33	0.59
1:C:33:SER:CA	1:C:36:VAL:HG23	2.29	0.59
1:C:139:VAL:HG22	1:C:146:GLN:HG2	1.83	0.59
1:D:96:LEU:O	1:D:98:VAL:N	2.35	0.59
1:A:11:ARG:HH11	1:A:11:ARG:HB2	1.68	0.59
1:C:181:LEU:CD1	1:C:184:MET:HE3	2.33	0.59
1:B:195:VAL:HG13	1:B:196:THR:H	1.69	0.58
1:B:5:LEU:C	1:B:7:ALA:H	2.11	0.58
1:D:110:ALA:HA	1:D:114:VAL:CG2	2.33	0.58
1:D:102:VAL:HG23	1:D:103:LEU:H	1.68	0.58
1:D:197:ARG:HG3	1:D:197:ARG:NH1	2.18	0.58
1:A:86:LEU:HD23	1:A:86:LEU:O	2.03	0.58
1:D:86:LEU:CD2	1:D:135:LEU:CD1	2.81	0.58
1:C:164:THR:O	1:C:167:VAL:HG12	2.03	0.58
1:C:163:GLY:O	1:C:167:VAL:HB	2.04	0.58
1:C:96:LEU:CB	1:C:103:LEU:HD12	2.33	0.58
1:B:14:ILE:HG21	1:B:50:PHE:CE2	2.39	0.58
3:F:16:DC:H2'	3:F:17:DG:C8	2.39	0.57
1:A:104:ARG:HG3	1:A:104:ARG:HH11	1.69	0.57
1:B:46:LEU:HD23	1:B:46:LEU:O	2.03	0.57
1:C:32:LEU:O	1:C:36:VAL:HG22	2.04	0.57
1:D:87:MET:O	1:D:90:THR:HB	2.05	0.57
1:A:28:GLU:C	1:A:30:THR:N	2.62	0.57
1:A:188:LEU:O	1:A:192:MET:HG2	2.05	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:191:GLY:HA2	1:D:190:ARG:HD3	1.86	0.57
1:B:64:GLN:HA	1:B:67:THR:HG22	1.87	0.57
1:C:137:ASP:O	1:C:140:ARG:HB2	2.03	0.57
1:D:44:GLY:C	3:F:1:DA:H8	2.11	0.56
1:D:96:LEU:C	1:D:98:VAL:H	2.12	0.56
1:B:86:LEU:HB2	1:B:134:ARG:CG	2.27	0.56
1:D:157:LEU:HD23	1:D:184:MET:HE1	1.86	0.56
1:A:151:ASP:HA	1:A:152:SER:CB	2.35	0.56
1:A:159:CYS:O	1:A:161:VAL:N	2.38	0.56
1:D:45:ALA:CB	3:F:1:DA:H2'	2.35	0.56
1:A:156:THR:HB	1:A:157:LEU:HG	1.87	0.56
1:C:198:ARG:O	1:C:202:VAL:HG23	2.05	0.56
1:B:7:ALA:CB	3:F:4:DA:H5''	2.34	0.55
1:B:204:LEU:O	1:B:208:LEU:CB	2.54	0.55
1:D:183:GLU:HA	1:D:186:TYR:CD2	2.40	0.55
1:A:69:ARG:HG2	1:A:69:ARG:NH1	2.22	0.55
1:A:34:GLU:HA	1:A:37:ALA:HB3	1.89	0.55
1:A:204:LEU:O	1:A:204:LEU:HD12	2.07	0.55
1:A:183:GLU:O	1:A:186:TYR:HB2	2.07	0.55
1:A:83:LEU:O	1:A:83:LEU:HD23	2.07	0.55
1:A:8:GLU:C	1:A:8:GLU:CD	2.75	0.55
1:A:100:GLY:HA3	1:A:103:LEU:HD12	1.88	0.55
1:A:153:VAL:O	1:A:156:THR:OG1	2.23	0.55
1:D:101:PRO:HG2	1:D:102:VAL:H	1.72	0.55
1:C:36:VAL:HG12	1:C:41:VAL:HG23	1.88	0.54
1:C:76:ASP:C	1:C:78:ARG:N	2.63	0.54
1:D:98:VAL:O	1:D:100:GLY:N	2.40	0.54
1:D:132:THR:HG23	1:D:136:LEU:HD23	1.88	0.54
1:D:96:LEU:C	1:D:98:VAL:N	2.63	0.54
1:B:175:GLY:H	1:B:177:GLU:HG3	1.72	0.54
1:D:181:LEU:O	1:D:182:ALA:C	2.49	0.54
1:C:139:VAL:HG23	1:C:144:VAL:HB	1.89	0.54
1:D:45:ALA:N	3:F:1:DA:H2'	2.23	0.54
1:A:6:ARG:NE	1:A:6:ARG:H	2.06	0.54
1:A:31:THR:CG2	1:A:34:GLU:HG2	2.38	0.54
1:B:61:LEU:O	1:B:64:GLN:HB3	2.08	0.54
1:C:208:LEU:O	1:C:212:THR:HB	2.08	0.54
3:F:9:DG:H2''	3:F:10:DG:H8	1.71	0.53
2:E:21:DA:N1	3:F:2:DT:O4	2.41	0.53
1:B:19:ALA:HB1	1:B:102:VAL:HA	1.90	0.53
2:E:16:DC:H2''	2:E:17:DG:N7	2.24	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:81:SER:O	1:B:82:SER:C	2.51	0.53
1:C:103:LEU:O	1:C:104:ARG:C	2.52	0.53
1:B:163:GLY:HA3	1:B:180:ARG:CZ	2.38	0.53
1:A:153:VAL:C	1:A:155:HIS:HD2	2.15	0.53
1:D:86:LEU:HD13	1:D:134:ARG:HG3	1.91	0.53
1:A:128:ARG:HG3	1:A:153:VAL:HG21	1.91	0.53
2:E:20:DT:H2''	2:E:21:DA:C8	2.44	0.53
1:B:21:LEU:CD1	1:B:25:ARG:NH2	2.72	0.52
2:E:9:DG:H2''	2:E:10:DA:H8	1.74	0.52
1:D:90:THR:HG22	1:D:91:PHE:N	2.25	0.52
1:B:30:THR:O	1:B:53:LYS:NZ	2.43	0.52
2:E:8:DG:H2''	2:E:9:DG:C8	2.45	0.52
1:A:28:GLU:HA	1:A:53:LYS:NZ	2.25	0.52
1:A:158:VAL:O	1:A:162:VAL:HG23	2.10	0.52
1:A:198:ARG:O	1:A:199:ALA:C	2.52	0.52
1:C:86:LEU:HD21	1:C:185:TRP:HH2	1.74	0.52
1:A:121:PRO:O	1:A:122:HIS:C	2.52	0.52
1:D:32:LEU:HD12	1:D:35:ILE:HD12	1.92	0.52
1:A:64:GLN:O	1:A:67:THR:N	2.43	0.52
1:C:95:ARG:HH11	1:C:96:LEU:HD11	1.75	0.52
1:D:137:ASP:CG	1:D:140:ARG:HH11	2.18	0.52
1:B:80:TYR:CG	1:B:84:GLU:HB2	2.44	0.51
1:B:139:VAL:HA	1:B:144:VAL:HG22	1.92	0.51
2:E:4:DT:H5'	2:E:4:DT:C6	2.45	0.51
2:E:19:DT:H2'	2:E:20:DT:H71	1.92	0.51
1:D:157:LEU:O	1:D:161:VAL:HG12	2.10	0.51
1:A:188:LEU:HB3	1:A:192:MET:HE2	1.92	0.51
1:A:202:VAL:O	1:A:203:THR:C	2.53	0.51
1:C:175:GLY:CA	1:C:176:ARG:HB3	2.29	0.51
1:D:43:LYS:O	1:D:46:LEU:HB3	2.10	0.51
2:E:4:DT:O4	3:F:19:DA:N1	2.44	0.51
2:E:6:DC:H2'	2:E:7:DG:C8	2.45	0.51
3:F:14:DC:OP2	3:F:14:DC:H2'	2.10	0.51
1:A:36:VAL:HB	1:A:41:VAL:O	2.10	0.51
1:B:136:LEU:O	1:B:136:LEU:HG	2.10	0.51
1:B:189:ILE:HG23	1:B:193:VAL:HG21	1.92	0.51
1:D:32:LEU:O	1:D:33:SER:C	2.50	0.51
3:F:2:DT:H2''	3:F:3:DA:C8	2.45	0.51
1:A:67:THR:HA	1:A:70:ARG:HB3	1.91	0.51
1:A:185:TRP:O	1:A:189:ILE:HG13	2.10	0.51
1:B:153:VAL:HG11	1:B:192:MET:SD	2.51	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:177:GLU:N	1:C:178:PRO:HD2	2.26	0.51
1:D:50:PHE:N	1:D:50:PHE:CD2	2.78	0.51
1:A:41:VAL:CG1	1:A:45:ALA:HB3	2.40	0.51
1:B:124:PHE:C	1:B:126:GLU:N	2.69	0.51
1:C:5:LEU:CD1	3:F:21:DG:OP1	2.59	0.51
1:C:56:LEU:O	1:C:57:ALA:C	2.54	0.51
1:C:47:TYR:CD1	3:F:13:DT:C7	2.93	0.51
1:C:164:THR:HG21	1:D:158:VAL:HG12	1.92	0.51
1:D:188:LEU:HB3	1:D:192:MET:CE	2.41	0.51
3:F:2:DT:C2'	3:F:3:DA:H5'	2.34	0.51
1:A:45:ALA:HA	2:E:3:DA:H3'	1.93	0.51
1:B:86:LEU:HD12	1:B:185:TRP:CH2	2.46	0.51
1:D:101:PRO:O	1:D:102:VAL:C	2.54	0.51
1:B:80:TYR:CG	1:B:84:GLU:CB	2.94	0.50
1:C:144:VAL:O	1:C:194:PRO:HD3	2.11	0.50
1:D:23:ASP:HB2	1:D:108:ARG:HG3	1.92	0.50
2:E:21:DA:C5	2:E:22:DT:H2'	2.45	0.50
2:E:7:DG:H1	3:F:16:DC:N4	2.04	0.50
2:E:17:DG:N2	3:F:6:DC:C2	2.72	0.50
1:A:71:LEU:HA	1:A:74:ASP:OD1	2.12	0.50
1:C:70:ARG:NH1	1:C:71:LEU:CD1	2.74	0.50
1:D:41:VAL:HG23	1:D:42:THR:O	2.11	0.50
1:D:45:ALA:HB2	3:F:1:DA:C2'	2.42	0.50
1:D:47:TYR:OH	2:E:17:DG:H2''	2.12	0.50
2:E:7:DG:N2	3:F:16:DC:C2	2.77	0.50
1:C:96:LEU:O	1:C:97:CYS:C	2.52	0.50
1:D:10:THR:O	1:D:14:ILE:HG13	2.12	0.49
1:D:68:SER:OG	1:D:93:MET:HE1	2.12	0.49
1:A:28:GLU:C	1:A:30:THR:H	2.20	0.49
1:C:135:LEU:O	1:C:138:ALA:N	2.45	0.49
1:D:45:ALA:HB2	3:F:1:DA:C3'	2.42	0.49
3:F:2:DT:H4'	3:F:2:DT:OP1	2.12	0.49
1:A:139:VAL:CG2	1:A:144:VAL:HB	2.42	0.49
1:C:5:LEU:N	1:C:8:GLU:H	2.10	0.49
1:A:150:VAL:C	1:A:152:SER:HB3	2.36	0.49
1:A:157:LEU:HA	1:A:184:MET:HE1	1.95	0.49
1:B:135:LEU:C	1:B:137:ASP:H	2.21	0.49
1:C:65:SER:OG	1:C:121:PRO:HB3	2.13	0.49
1:A:186:TYR:CD1	1:A:202:VAL:HG13	2.48	0.48
3:F:1:DA:H1'	3:F:2:DT:P	2.52	0.48
2:E:6:DC:C2'	2:E:7:DG:H8	2.19	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:25:ARG:HG3	1:D:25:ARG:NH2	2.28	0.48
1:D:117:ARG:HE	1:D:120:LEU:N	2.10	0.48
1:D:141:GLN:O	1:D:142:SER:HB2	2.12	0.48
1:A:151:ASP:HB3	1:A:153:VAL:CG2	2.30	0.48
1:B:10:THR:O	1:B:11:ARG:C	2.57	0.48
1:A:196:THR:CG2	1:A:197:ARG:HD2	2.43	0.48
1:B:139:VAL:CG2	1:B:144:VAL:HG21	2.39	0.48
1:B:175:GLY:N	1:B:177:GLU:OE1	2.43	0.48
1:D:86:LEU:HD21	1:D:135:LEU:CD1	2.38	0.48
1:D:157:LEU:CD2	1:D:184:MET:HE1	2.44	0.48
1:B:32:LEU:HD21	1:B:53:LYS:HB2	1.96	0.48
1:B:49:HIS:NE2	3:F:5:DA:OP2	2.38	0.48
1:A:190:ARG:NH2	1:B:145:HIS:CD2	2.82	0.47
1:A:65:SER:O	1:A:69:ARG:HG3	2.14	0.47
1:A:159:CYS:C	1:A:161:VAL:N	2.73	0.47
1:C:92:GLY:O	1:C:96:LEU:HD22	2.14	0.47
1:A:6:ARG:H	1:A:6:ARG:CZ	2.27	0.47
1:A:47:TYR:OH	3:F:16:DC:OP2	2.21	0.47
1:C:176:ARG:O	1:C:179:ARG:HB2	2.14	0.47
1:D:160:SER:C	1:D:162:VAL:H	2.23	0.47
3:F:9:DG:H8	3:F:9:DG:H5''	1.79	0.47
1:B:195:VAL:HG13	1:B:196:THR:N	2.30	0.47
1:D:188:LEU:HB3	1:D:192:MET:HE2	1.95	0.47
1:A:28:GLU:H	1:A:28:GLU:CD	2.19	0.47
1:B:33:SER:HA	1:B:36:VAL:HG22	1.97	0.47
1:D:42:THR:CG2	3:F:2:DT:OP2	2.61	0.47
1:D:72:ALA:O	1:D:75:LEU:HG	2.14	0.47
1:D:189:ILE:HD12	1:D:201:TYR:HB3	1.97	0.47
1:A:36:VAL:HG23	1:A:37:ALA:N	2.29	0.47
1:A:78:ARG:HD2	1:A:80:TYR:CE2	2.50	0.47
1:C:180:ARG:HA	1:C:183:GLU:OE2	2.14	0.47
1:D:105:ALA:O	1:D:109:LEU:HG	2.15	0.47
2:E:18:DT:C2'	2:E:19:DT:C6	2.97	0.47
1:C:6:ARG:NH2	2:E:6:DC:H4'	2.29	0.47
1:D:194:PRO:O	1:D:198:ARG:HB2	2.15	0.47
1:A:28:GLU:O	1:A:29:SER:C	2.58	0.47
1:A:159:CYS:HA	1:A:162:VAL:HB	1.97	0.47
1:B:163:GLY:HA3	1:B:180:ARG:HH22	1.78	0.47
1:B:199:ALA:O	1:B:201:TYR:N	2.48	0.47
1:A:185:TRP:O	1:A:186:TYR:C	2.58	0.47
1:C:163:GLY:C	1:C:166:VAL:HB	2.40	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:190:ARG:HE	1:B:198:ARG:NH2	2.03	0.46
1:B:44:GLY:HA3	3:F:7:DG:O6	2.15	0.46
1:B:165:ARG:O	1:B:171:LEU:N	2.42	0.46
1:C:47:TYR:CE1	1:C:51:ALA:HA	2.50	0.46
1:C:157:LEU:O	1:C:161:VAL:HG23	2.16	0.46
1:C:76:ASP:O	1:C:78:ARG:N	2.49	0.46
1:C:156:THR:HG22	1:C:160:SER:OG	2.16	0.46
1:C:199:ALA:O	1:C:200:ARG:C	2.59	0.46
1:B:201:TYR:HA	1:B:204:LEU:HB3	1.97	0.46
1:C:189:ILE:HD13	1:C:201:TYR:HB3	1.98	0.46
1:C:117:ARG:H	1:C:120:LEU:HD13	1.81	0.46
1:D:69:ARG:HH11	1:D:69:ARG:CB	2.27	0.46
1:B:116:VAL:O	1:B:117:ARG:CB	2.64	0.46
1:D:44:GLY:O	3:F:1:DA:C8	2.66	0.46
1:B:96:LEU:O	1:B:97:CYS:C	2.59	0.45
1:C:109:LEU:HD23	1:C:109:LEU:HA	1.75	0.45
2:E:21:DA:C4	2:E:22:DT:H2'	2.52	0.45
1:C:148:ILE:HD13	1:C:148:ILE:HA	1.77	0.45
1:D:148:ILE:HG21	1:D:192:MET:HG2	1.98	0.45
1:A:107:LEU:O	1:A:111:THR:HG23	2.15	0.45
1:C:78:ARG:HD2	1:C:80:TYR:HE2	1.81	0.45
1:C:82:SER:HB2	1:C:141:GLN:NE2	2.31	0.45
1:D:136:LEU:HD21	1:D:150:VAL:HG21	1.98	0.45
1:A:86:LEU:CD1	1:A:135:LEU:CD1	2.91	0.45
1:D:37:ALA:C	1:D:39:ALA:H	2.25	0.45
2:E:21:DA:C8	2:E:22:DT:H2'	2.51	0.45
3:F:14:DC:H2''	3:F:15:DC:C6	2.51	0.45
1:A:159:CYS:O	1:A:162:VAL:N	2.50	0.45
1:B:19:ALA:CB	1:B:60:ILE:HD11	2.46	0.45
1:D:197:ARG:HH11	1:D:197:ARG:CG	2.29	0.45
1:A:28:GLU:O	1:A:30:THR:N	2.50	0.45
1:D:96:LEU:O	1:D:99:GLN:N	2.50	0.45
1:D:178:PRO:C	1:D:180:ARG:H	2.24	0.45
1:D:42:THR:CB	3:F:2:DT:OP2	2.65	0.45
1:A:153:VAL:C	1:A:155:HIS:CD2	2.95	0.45
1:D:84:GLU:O	1:D:88:ARG:HG3	2.17	0.45
1:B:179:ARG:NH2	1:B:179:ARG:HB3	2.32	0.45
1:D:159:CYS:O	1:D:161:VAL:N	2.50	0.45
1:B:100:GLY:HA2	1:B:101:PRO:HD2	1.84	0.44
1:D:62:GLU:O	1:D:66:ARG:HB2	2.17	0.44
1:D:149:ASP:O	1:D:153:VAL:HG23	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:46:LEU:HD23	1:C:50:PHE:HB2	1.99	0.44
1:B:135:LEU:C	1:B:137:ASP:N	2.76	0.44
1:C:179:ARG:O	1:C:182:ALA:HB3	2.17	0.44
1:C:70:ARG:NH1	1:C:71:LEU:HD11	2.32	0.44
1:C:17:ALA:CB	1:C:39:ALA:HB2	2.46	0.44
1:D:42:THR:HG21	3:F:1:DA:C3'	2.47	0.44
1:D:45:ALA:HB2	3:F:1:DA:H3'	1.98	0.44
1:B:46:LEU:HD23	1:B:46:LEU:C	2.42	0.44
1:B:62:GLU:O	1:B:64:GLN:N	2.51	0.44
1:A:8:GLU:O	1:A:11:ARG:N	2.51	0.44
1:A:71:LEU:O	1:A:71:LEU:HG	2.17	0.44
1:A:73:LYS:HA	1:A:73:LYS:HD3	1.61	0.44
1:A:104:ARG:HG3	1:A:104:ARG:NH1	2.32	0.44
1:B:17:ALA:CB	1:B:39:ALA:HB2	2.48	0.44
1:B:201:TYR:O	1:B:204:LEU:N	2.51	0.44
1:C:6:ARG:HE	1:C:6:ARG:HB3	1.62	0.44
1:A:10:THR:O	1:A:13:THR:N	2.51	0.44
1:A:41:VAL:HG12	1:A:42:THR:O	2.17	0.44
1:A:123:PRO:O	1:A:126:GLU:HB3	2.18	0.44
1:B:11:ARG:C	1:B:11:ARG:HD3	2.43	0.44
1:C:31:THR:C	1:C:33:SER:N	2.75	0.44
1:C:135:LEU:C	1:C:137:ASP:N	2.76	0.44
1:C:33:SER:HB3	1:C:43:LYS:HE2	1.98	0.43
1:C:139:VAL:HG13	1:C:140:ARG:HH22	1.83	0.43
1:B:124:PHE:O	1:B:127:TRP:N	2.50	0.43
1:C:73:LYS:C	1:C:75:LEU:H	2.26	0.43
1:D:48:PHE:HD2	1:D:49:HIS:HD2	1.65	0.43
1:D:130:ILE:O	1:D:131:ALA:C	2.61	0.43
1:D:177:GLU:HB2	1:D:178:PRO:CD	2.47	0.43
1:C:87:MET:HE2	1:C:204:LEU:HG	1.98	0.43
1:D:48:PHE:CD2	1:D:49:HIS:HD2	2.36	0.43
1:A:190:ARG:HH22	1:B:145:HIS:CD2	2.36	0.43
1:B:44:GLY:CA	3:F:7:DG:O6	2.66	0.43
1:C:138:ALA:O	1:C:142:SER:O	2.35	0.43
1:D:41:VAL:HG23	1:D:45:ALA:HB3	2.00	0.43
1:C:82:SER:HB3	1:C:138:ALA:HB2	1.99	0.43
1:C:141:GLN:H	1:C:141:GLN:HG2	1.47	0.43
1:A:23:ASP:HA	1:A:108:ARG:NE	2.33	0.43
1:B:185:TRP:O	1:B:186:TYR:C	2.60	0.43
1:A:73:LYS:HD3	1:A:76:ASP:OD2	2.18	0.43
1:A:139:VAL:HG23	1:A:144:VAL:HB	1.99	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:F:22:DT:O2	3:F:22:DT:O4'	2.34	0.43
1:C:185:TRP:O	1:C:186:TYR:C	2.61	0.43
1:B:32:LEU:HD21	1:B:53:LYS:CB	2.49	0.42
1:C:139:VAL:HG12	1:C:140:ARG:HH12	1.84	0.42
3:F:15:DC:H2''	3:F:16:DC:OP2	2.18	0.42
1:A:36:VAL:HG11	1:A:43:LYS:HA	1.99	0.42
1:A:41:VAL:HG13	1:A:45:ALA:HB3	2.00	0.42
1:A:149:ASP:O	1:A:151:ASP:N	2.52	0.42
1:B:27:TYR:OH	1:B:53:LYS:HG3	2.19	0.42
1:B:103:LEU:HD23	1:B:103:LEU:HA	1.60	0.42
1:C:15:ILE:O	1:C:16:GLY:C	2.61	0.42
1:D:10:THR:OG1	1:D:49:HIS:CE1	2.72	0.42
2:E:7:DG:H2''	2:E:8:DG:OP2	2.20	0.42
1:B:113:GLY:O	1:B:115:PRO:HD3	2.19	0.42
1:B:177:GLU:HB2	1:B:178:PRO:HD3	2.02	0.42
1:A:28:GLU:HA	1:A:53:LYS:HZ3	1.84	0.42
1:A:32:LEU:O	1:A:36:VAL:HG22	2.20	0.42
1:B:81:SER:H	1:B:84:GLU:CD	2.27	0.42
1:C:10:THR:O	1:C:11:ARG:C	2.63	0.42
1:C:96:LEU:HA	1:C:99:GLN:HG2	2.01	0.42
1:D:10:THR:OG1	1:D:49:HIS:HE1	2.03	0.42
1:B:81:SER:O	1:B:84:GLU:OE2	2.37	0.42
1:B:145:HIS:CE1	1:B:194:PRO:HD3	2.55	0.42
1:D:116:VAL:HG22	1:D:120:LEU:HB3	2.00	0.42
1:A:125:THR:HG22	1:A:128:ARG:NH2	2.33	0.42
1:B:83:LEU:O	1:B:87:MET:HB2	2.19	0.42
1:D:159:CYS:O	1:D:160:SER:C	2.62	0.42
1:D:178:PRO:C	1:D:180:ARG:N	2.77	0.42
3:F:2:DT:C2'	3:F:3:DA:C8	3.02	0.42
1:A:82:SER:H	1:A:141:GLN:HE22	1.67	0.42
1:C:87:MET:HE3	1:C:205:ALA:HB2	2.01	0.42
1:D:47:TYR:CE1	2:E:18:DT:H5'	2.54	0.42
1:B:14:ILE:CG2	1:B:50:PHE:HE2	2.30	0.42
1:C:19:ALA:CB	1:C:60:ILE:HD11	2.50	0.42
1:B:143:ASP:O	1:B:194:PRO:HD2	2.20	0.42
1:C:84:GLU:HG3	1:C:204:LEU:CD2	2.50	0.42
1:D:14:ILE:HD13	1:D:46:LEU:HA	2.02	0.42
1:D:45:ALA:HB2	3:F:1:DA:H2'	2.01	0.42
1:D:83:LEU:HD12	1:D:201:TYR:CD1	2.55	0.42
1:A:119:PRO:O	1:A:120:LEU:C	2.63	0.41
1:C:31:THR:C	1:C:33:SER:H	2.26	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:85:ALA:O	1:D:89:LEU:N	2.49	0.41
1:A:158:VAL:HG12	1:A:159:CYS:N	2.35	0.41
1:A:75:LEU:C	1:A:77:GLY:H	2.28	0.41
1:B:91:PHE:HE2	1:B:185:TRP:CD1	2.39	0.41
1:B:118:PRO:HA	1:B:119:PRO:HA	1.82	0.41
1:C:44:GLY:HA3	2:E:9:DG:O6	2.21	0.41
1:D:178:PRO:HB2	1:D:209:GLU:HA	2.03	0.41
2:E:11:DC:C2'	2:E:12:DG:N7	2.82	0.41
1:A:60:ILE:HA	1:A:63:ILE:HD12	2.03	0.41
1:C:13:THR:HG22	1:C:39:ALA:HB1	2.02	0.41
1:C:73:LYS:O	1:C:75:LEU:N	2.54	0.41
1:C:96:LEU:N	1:C:96:LEU:HD13	2.34	0.41
3:F:9:DG:H5''	3:F:9:DG:C8	2.56	0.41
3:F:16:DC:H2'	3:F:17:DG:N7	2.36	0.41
1:A:89:LEU:C	1:A:91:PHE:N	2.78	0.41
1:B:54:GLU:CG	1:B:115:PRO:HB2	2.50	0.41
1:C:32:LEU:HB2	3:F:11:DC:OP2	2.20	0.41
1:B:62:GLU:C	1:B:64:GLN:N	2.74	0.41
1:B:145:HIS:HB2	1:B:148:ILE:HG23	2.02	0.41
3:F:21:DG:N2	3:F:22:DT:O4	2.54	0.41
1:A:130:ILE:CG2	1:A:131:ALA:N	2.84	0.41
1:B:138:ALA:HB1	1:B:143:ASP:HB2	2.03	0.41
1:B:139:VAL:CA	1:B:144:VAL:HG22	2.50	0.41
1:B:179:ARG:NH1	1:B:209:GLU:OE1	2.54	0.41
1:C:6:ARG:HD2	2:E:7:DG:H5'	2.03	0.41
1:C:19:ALA:HB1	1:C:102:VAL:HA	2.03	0.41
1:D:180:ARG:NH1	1:D:180:ARG:CG	2.72	0.41
3:F:1:DA:H2''	3:F:2:DT:OP2	2.10	0.41
1:A:193:VAL:O	1:A:194:PRO:C	2.62	0.41
1:C:28:GLU:CG	1:D:28:GLU:HB3	2.50	0.41
1:C:42:THR:O	1:C:45:ALA:N	2.54	0.41
1:C:164:THR:C	1:C:167:VAL:H	2.28	0.41
1:C:208:LEU:HD23	1:C:208:LEU:HA	1.79	0.41
1:D:101:PRO:O	1:D:104:ARG:N	2.54	0.41
1:B:175:GLY:H	1:B:177:GLU:CG	2.33	0.40
1:B:176:ARG:NH1	1:B:179:ARG:HD3	2.36	0.40
1:B:189:ILE:HG23	1:B:193:VAL:CG2	2.50	0.40
1:D:66:ARG:O	1:D:69:ARG:HG2	2.21	0.40
1:B:5:LEU:C	1:B:7:ALA:N	2.76	0.40
1:A:150:VAL:HG23	1:A:151:ASP:OD2	2.22	0.40
1:B:21:LEU:HD11	1:B:25:ARG:HH21	1.84	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:127:TRP:O	1:B:128:ARG:C	2.65	0.40
1:C:73:LYS:HD3	1:C:74:ASP:N	2.36	0.40
1:C:165:ARG:HG2	1:C:165:ARG:O	2.20	0.40
1:B:80:TYR:CG	1:B:84:GLU:HB3	2.57	0.40
1:B:122:HIS:HA	1:B:123:PRO:HD3	1.86	0.40
1:B:204:LEU:O	1:B:208:LEU:HD12	2.22	0.40
1:C:5:LEU:HB3	3:F:21:DG:H5'	2.03	0.40
1:D:32:LEU:HA	1:D:32:LEU:HD12	1.90	0.40
1:D:136:LEU:HA	1:D:136:LEU:HD13	1.50	0.40
1:C:6:ARG:CZ	2:E:6:DC:H4'	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	
1	A	188/215 (87%)	154 (82%)	19 (10%)	15 (8%)	1	4
1	B	200/215 (93%)	145 (72%)	41 (20%)	14 (7%)	1	6
1	C	194/215 (90%)	136 (70%)	45 (23%)	13 (7%)	1	7
1	D	185/215 (86%)	145 (78%)	30 (16%)	10 (5%)	1	12
All	All	767/860 (89%)	580 (76%)	135 (18%)	52 (7%)	1	7

All (52) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	70	ARG
1	A	153	VAL
1	A	160	SER
1	A	177	GLU
1	A	210	GLN

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Mol	Chain	Res	Type
1	B	78	ARG
1	B	80	TYR
1	B	117	ARG
1	B	142	SER
1	B	190	ARG
1	C	162	VAL
1	D	99	GLN
1	D	102	VAL
1	D	121	PRO
1	A	29	SER
1	A	74	ASP
1	A	152	SER
1	A	154	ALA
1	B	6	ARG
1	B	208	LEU
1	C	77	GLY
1	C	198	ARG
1	D	97	CYS
1	A	39	ALA
1	A	90	THR
1	A	199	ALA
1	B	9	GLN
1	B	82	SER
1	B	136	LEU
1	B	200	ARG
1	C	85	ALA
1	C	86	LEU
1	C	144	VAL
1	C	199	ALA
1	D	101	PRO
1	D	161	VAL
1	D	177	GLU
1	B	63	ILE
1	C	136	LEU
1	D	38	HIS
1	A	69	ARG
1	A	78	ARG
1	B	112	ALA
1	B	173	PRO
1	C	74	ASP
1	C	122	HIS
1	D	66	ARG

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Mol	Chain	Res	Type
1	D	160	SER
1	A	150	VAL
1	C	16	GLY
1	C	41	VAL
1	C	163	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	
1	A	148/171 (86%)	104 (70%)	44 (30%)	0	1
1	B	148/171 (86%)	119 (80%)	29 (20%)	1	8
1	C	153/171 (90%)	114 (74%)	39 (26%)	0	3
1	D	154/171 (90%)	118 (77%)	36 (23%)	1	4
All	All	603/684 (88%)	455 (76%)	148 (24%)	1	3

All (148) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	ARG
1	A	11	ARG
1	A	21	LEU
1	A	28	GLU
1	A	29	SER
1	A	32	LEU
1	A	33	SER
1	A	42	THR
1	A	46	LEU
1	A	56	LEU
1	A	62	GLU
1	A	65	SER
1	A	67	THR
1	A	69	ARG
1	A	73	LYS
1	A	74	ASP

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Mol	Chain	Res	Type
1	A	76	ASP
1	A	86	LEU
1	A	99	GLN
1	A	102	VAL
1	A	103	LEU
1	A	108	ARG
1	A	111	THR
1	A	130	ILE
1	A	135	LEU
1	A	136	LEU
1	A	142	SER
1	A	146	GLN
1	A	148	ILE
1	A	151	ASP
1	A	155	HIS
1	A	156	THR
1	A	157	LEU
1	A	158	VAL
1	A	159	CYS
1	A	160	SER
1	A	177	GLU
1	A	180	ARG
1	A	181	LEU
1	A	184	MET
1	A	196	THR
1	A	204	LEU
1	A	207	ARG
1	A	208	LEU
1	B	6	ARG
1	B	8	GLU
1	B	15	ILE
1	B	21	LEU
1	B	31	THR
1	B	60	ILE
1	B	61	LEU
1	B	65	SER
1	B	75	LEU
1	B	84	GLU
1	B	86	LEU
1	B	98	VAL
1	B	99	GLN
1	B	111	THR

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Mol	Chain	Res	Type
1	B	125	THR
1	B	134	ARG
1	B	136	LEU
1	B	147	ASP
1	B	150	VAL
1	B	152	SER
1	B	153	VAL
1	B	160	SER
1	B	179	ARG
1	B	187	ILE
1	B	195	VAL
1	B	197	ARG
1	B	208	LEU
1	B	210	GLN
1	B	212	THR
1	C	6	ARG
1	C	15	ILE
1	C	29	SER
1	C	30	THR
1	C	31	THR
1	C	32	LEU
1	C	36	VAL
1	C	54	GLU
1	C	60	ILE
1	C	62	GLU
1	C	71	LEU
1	C	73	LYS
1	C	75	LEU
1	C	82	SER
1	C	86	LEU
1	C	96	LEU
1	C	102	VAL
1	C	107	LEU
1	C	120	LEU
1	C	125	THR
1	C	133	SER
1	C	134	ARG
1	C	140	ARG
1	C	141	GLN
1	C	143	ASP
1	C	148	ILE
1	C	157	LEU

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Mol	Chain	Res	Type
1	C	158	VAL
1	C	162	VAL
1	C	164	THR
1	C	167	VAL
1	C	177	GLU
1	C	179	ARG
1	C	188	LEU
1	C	195	VAL
1	C	196	THR
1	C	209	GLU
1	C	211	GLU
1	C	212	THR
1	D	8	GLU
1	D	23	ASP
1	D	32	LEU
1	D	41	VAL
1	D	42	THR
1	D	54	GLU
1	D	73	LYS
1	D	80	TYR
1	D	89	LEU
1	D	99	GLN
1	D	102	VAL
1	D	107	LEU
1	D	116	VAL
1	D	120	LEU
1	D	122	HIS
1	D	129	GLU
1	D	132	THR
1	D	133	SER
1	D	135	LEU
1	D	136	LEU
1	D	146	GLN
1	D	151	ASP
1	D	156	THR
1	D	157	LEU
1	D	158	VAL
1	D	164	THR
1	D	165	ARG
1	D	176	ARG
1	D	179	ARG
1	D	180	ARG

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Mol	Chain	Res	Type
1	D	184	MET
1	D	187	ILE
1	D	189	ILE
1	D	194	PRO
1	D	197	ARG
1	D	211	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	38	HIS
1	A	145	HIS
1	A	155	HIS
1	B	38	HIS
1	B	122	HIS
1	B	141	GLN
1	B	145	HIS
1	C	38	HIS
1	C	58	HIS
1	C	64	GLN
1	C	145	HIS
1	C	210	GLN
1	D	49	HIS
1	D	58	HIS
1	D	64	GLN
1	D	141	GLN

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

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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	A	194/215 (90%)	-1.09	0 100 100	42, 83, 130, 148	1 (0%)
1	B	204/215 (94%)	-1.01	0 100 100	51, 89, 143, 199	0
1	C	200/215 (93%)	-0.96	0 100 100	45, 89, 149, 186	1 (0%)
1	D	193/215 (89%)	-1.04	0 100 100	52, 82, 131, 169	0
2	E	20/22 (90%)	-0.93	0 100 100	52, 65, 102, 112	1 (5%)
3	F	22/22 (100%)	-0.89	0 100 100	53, 64, 119, 151	3 (13%)
All	All	833/904 (92%)	-1.02	0 100 100	42, 85, 138, 199	6 (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.