



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 27, 2023 – 11:26 PM EDT

PDB ID : 3LAJ
Title : The Structure of the Intermediate Complex of the Arginine Repressor from Mycobacterium tuberculosis Bound to its DNA Operator and L-arginine.
Authors : Cherney, L.T.; Cherney, M.M.; Garen, C.R.; James, M.N.G.; Mycobacterium Tuberculosis Structural Proteomics Project (XMTB)
Deposited on : 2010-01-06
Resolution : 2.31 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 : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.35
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.35

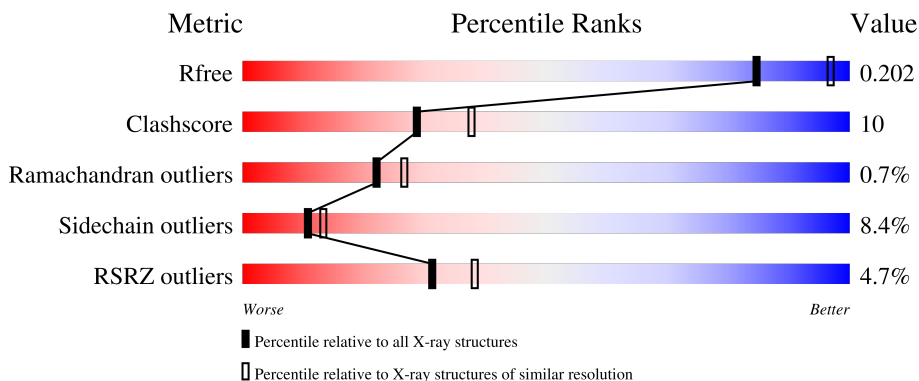
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.31 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

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	170	 2% 69% 16% 12%
1	B	170	 % 72% 13% 12%
1	C	170	 % 72% 13% 12%
1	D	170	 9% 65% 23% 9%
1	E	170	 8% 64% 24% 9%

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Mol	Chain	Length	Quality of chain
1	F	170	<p>6% 66% 21% 9%</p>
2	G	16	<p>38% 56% 6%</p>
2	I	16	<p>62% 31% 6%</p>
2	K	16	<p>38% 62%</p>
3	H	16	<p>44% 44% 12%</p>
3	J	16	<p>56% 31% 12%</p>
3	L	16	<p>38% 50% 12%</p>

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 9028 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 Arginine repressor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	149	1075	661	198	214	2	0	0	0
1	B	149	1075	661	198	214	2	0	0	0
1	C	149	1075	661	198	214	2	0	0	0
1	D	155	1117	685	207	223	2	0	0	0
1	E	154	1112	682	206	222	2	0	0	0
1	F	155	1117	685	207	223	2	0	0	0

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
2	G	16	326	157	62	92	15	0	0	0
2	I	16	326	157	62	92	15	0	0	0
2	K	16	326	157	62	92	15	0	0	0

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

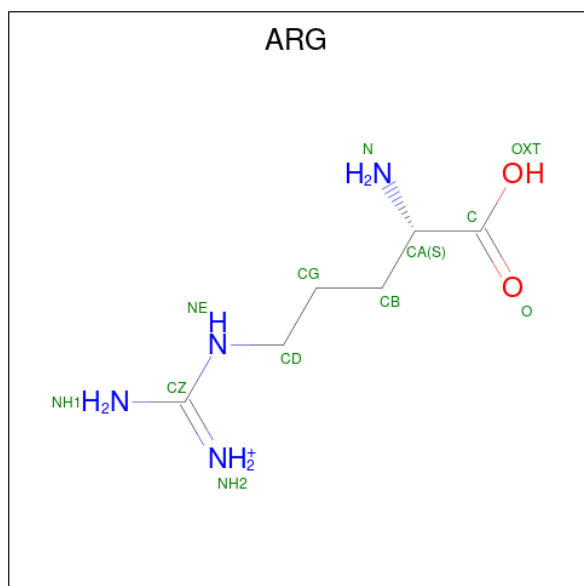
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
3	H	16	324	157	56	96	15	0	0	0
3	J	16	324	157	56	96	15	0	0	0

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
3	L	16	324	157	56	96	15	0	0	0

- Molecule 4 is ARGININE (three-letter code: ARG) (formula: C₆H₁₅N₄O₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
4	A	1	12	6	4	2	0	0
4	B	1	12	6	4	2	0	0
4	C	1	12	6	4	2	0	0
4	D	1	12	6	4	2	0	0
4	E	1	12	6	4	2	0	0
4	F	1	12	6	4	2	0	0

- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	64	Total 64 O 64	0	0
5	B	59	Total 59 O 59	0	0

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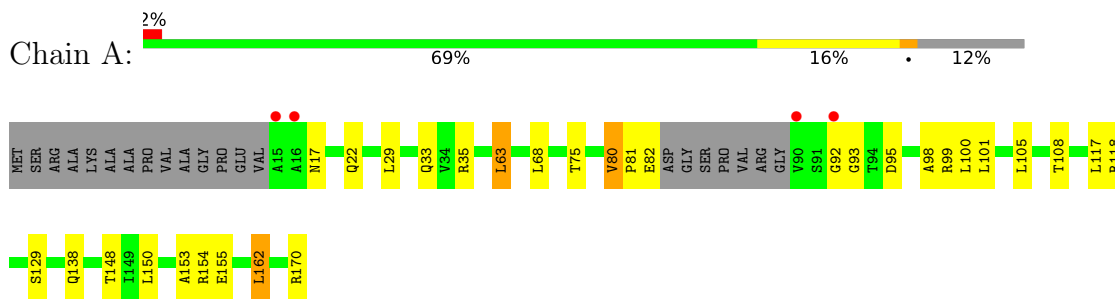
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	C	66	Total 66	O 66	0	0
5	D	47	Total 47	O 47	0	0
5	E	44	Total 44	O 44	0	0
5	F	45	Total 45	O 45	0	0
5	G	18	Total 18	O 18	0	0
5	H	25	Total 25	O 25	0	0
5	I	13	Total 13	O 13	0	0
5	J	20	Total 20	O 20	0	0
5	K	19	Total 19	O 19	0	0
5	L	15	Total 15	O 15	0	0

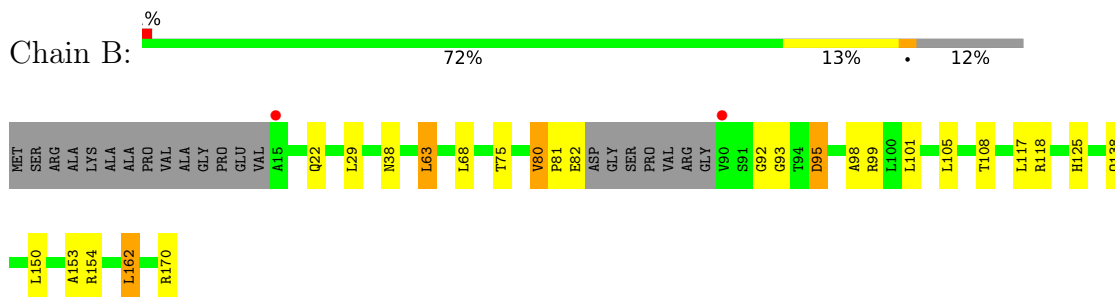
3 Residue-property plots [i](#)

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.

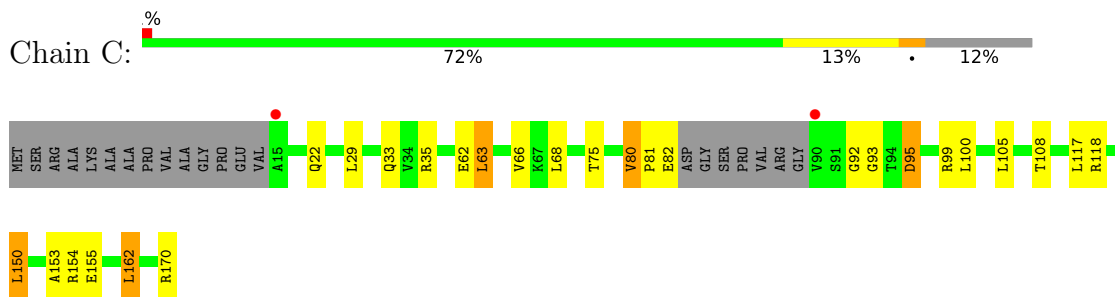
- Molecule 1: Arginine repressor



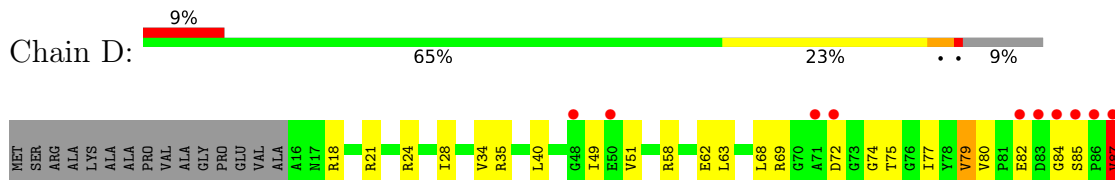
- Molecule 1: Arginine repressor

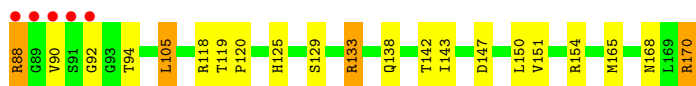


- Molecule 1: Arginine repressor

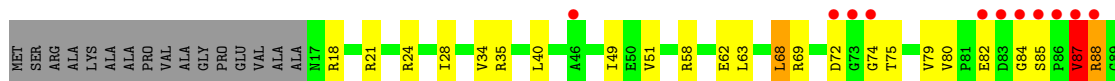


- Molecule 1: Arginine repressor

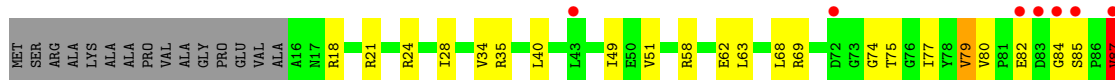




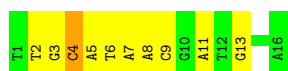
- Molecule 1: Arginine repressor



- Molecule 1: Arginine repressor



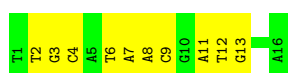
- Molecule 2: 5'-D(*TP*TP*GP*CP*AP*TP*AP*AP*CP*GP*AP*TP*GP*CP*AP*A)-3'



- Molecule 2: 5'-D(*TP*TP*GP*CP*AP*TP*AP*AP*CP*GP*AP*TP*GP*CP*AP*A)-3'



- Molecule 2: 5'-D(*TP*TP*GP*CP*AP*TP*AP*AP*CP*GP*AP*TP*GP*CP*AP*A)-3'



- Molecule 3: 5'-D(*TP*TP*GP*CP*AP*TP*CP*GP*TP*TP*AP*TP*GP*CP*AP*A)-3'





- Molecule 3: 5'-D(*TP*TP*GP*CP*AP*TP*CP*GP*TP*TP*AP*TP*GP*CP*AP*A)-3'



- Molecule 3: 5'-D(*TP*TP*GP*CP*AP*TP*CP*GP*TP*TP*AP*TP*GP*CP*AP*A)-3'



4 Data and refinement statistics i

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	185.69Å 106.79Å 119.08Å 90.00° 121.75° 90.00°	Depositor
Resolution (Å)	39.24 – 2.31 39.24 – 2.31	Depositor EDS
% Data completeness (in resolution range)	96.9 (39.24-2.31) 95.3 (39.24-2.31)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	0.04	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.13 (at 2.31Å)	Xtrriage
Refinement program	PHENIX (phenix.refine: 1.4_58)	Depositor
R, R_{free}	0.184 , 0.207 0.177 , 0.202	Depositor DCC
R_{free} test set	4221 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	50.7	Xtrriage
Anisotropy	0.268	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 49.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	0.417 for $1/2^*h-3/2^*k,-1/2^*h-1/2^*k,-1/2^*h+1/2^*k-1$ 0.419 for $1/2^*h+3/2^*k,1/2^*h-1/2^*k,-1/2^*h-1/2^*k-1$	Xtrriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	9028	wwPDB-VP
Average B, all atoms (Å ²)	58.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.67% 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	0.67	1/1082 (0.1%)	0.92	5/1467 (0.3%)
1	B	0.76	2/1082 (0.2%)	1.22	9/1467 (0.6%)
1	C	0.67	2/1082 (0.2%)	1.05	7/1467 (0.5%)
1	D	0.58	0/1126	0.91	6/1528 (0.4%)
1	E	0.69	2/1121 (0.2%)	1.15	8/1521 (0.5%)
1	F	0.61	1/1126 (0.1%)	1.01	7/1528 (0.5%)
2	G	1.09	0/366	1.78	8/563 (1.4%)
2	I	1.16	0/366	1.75	6/563 (1.1%)
2	K	1.21	0/366	1.77	8/563 (1.4%)
3	H	1.03	0/362	1.94	12/557 (2.2%)
3	J	1.02	0/362	1.84	9/557 (1.6%)
3	L	1.04	0/362	1.90	12/557 (2.2%)
All	All	0.79	8/8803 (0.1%)	1.31	97/12338 (0.8%)

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	154	ARG	CZ-NH1	-10.23	1.19	1.33
1	B	154	ARG	CZ-NH1	-9.89	1.20	1.33
1	B	154	ARG	CZ-NH2	-8.79	1.21	1.33
1	E	154	ARG	CZ-NH2	-8.04	1.22	1.33
1	C	154	ARG	CZ-NH2	-5.72	1.25	1.33
1	A	154	ARG	CZ-NH2	-5.69	1.25	1.33
1	F	154	ARG	CZ-NH2	-5.44	1.25	1.33
1	C	154	ARG	CZ-NH1	-5.06	1.26	1.33

All (97) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	170	ARG	NE-CZ-NH2	18.63	129.61	120.30
1	C	118	ARG	NE-CZ-NH1	-16.56	112.02	120.30
1	B	154	ARG	NE-CZ-NH2	16.01	128.30	120.30
1	C	118	ARG	NE-CZ-NH2	15.09	127.85	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	170	ARG	NE-CZ-NH1	-15.09	112.76	120.30
1	E	170	ARG	NE-CZ-NH2	14.31	127.46	120.30
1	E	170	ARG	NE-CZ-NH1	-14.29	113.16	120.30
1	F	133	ARG	NE-CZ-NH1	-14.23	113.19	120.30
1	E	154	ARG	NE-CZ-NH2	13.83	127.21	120.30
1	F	133	ARG	NE-CZ-NH2	13.10	126.85	120.30
3	L	11	DA	O4'-C1'-N9	-12.67	99.13	108.00
3	H	11	DA	O4'-C1'-N9	-12.18	99.47	108.00
1	E	154	ARG	NH1-CZ-NH2	-12.01	106.19	119.40
1	E	154	ARG	NE-CZ-NH1	11.97	126.29	120.30
1	B	154	ARG	NH1-CZ-NH2	-11.15	107.13	119.40
2	G	11	DA	O4'-C1'-N9	-10.93	100.35	108.00
2	I	11	DA	O4'-C1'-N9	-10.65	100.54	108.00
2	K	11	DA	O4'-C1'-N9	-10.16	100.89	108.00
1	B	118	ARG	NE-CZ-NH2	-10.02	115.29	120.30
1	A	118	ARG	NE-CZ-NH2	-9.86	115.37	120.30
3	L	3	DG	O4'-C1'-N9	-9.73	101.19	108.00
1	D	133	ARG	NE-CZ-NH2	-9.58	115.51	120.30
3	H	3	DG	O4'-C1'-N9	-9.47	101.37	108.00
1	E	133	ARG	NE-CZ-NH2	-9.41	115.59	120.30
1	C	154	ARG	NE-CZ-NH1	9.40	125.00	120.30
3	J	11	DA	O4'-C1'-N9	-9.29	101.50	108.00
1	B	170	ARG	CD-NE-CZ	9.19	136.46	123.60
2	I	3	DG	O4'-C1'-N9	-9.01	101.69	108.00
1	A	154	ARG	NE-CZ-NH1	8.95	124.77	120.30
3	J	3	DG	O4'-C1'-N9	-8.63	101.96	108.00
1	A	118	ARG	NE-CZ-NH1	8.56	124.58	120.30
1	B	118	ARG	NE-CZ-NH1	8.56	124.58	120.30
1	B	154	ARG	NE-CZ-NH1	8.53	124.56	120.30
1	E	133	ARG	NE-CZ-NH1	8.45	124.53	120.30
1	A	170	ARG	NE-CZ-NH1	8.28	124.44	120.30
1	C	170	ARG	NE-CZ-NH1	8.00	124.30	120.30
2	K	3	DG	O4'-C1'-N9	-7.99	102.41	108.00
1	C	118	ARG	CD-NE-CZ	7.84	134.58	123.60
2	G	3	DG	O4'-C1'-N9	-7.81	102.53	108.00
3	J	2	DT	O4'-C1'-N1	-7.80	102.54	108.00
1	D	170	ARG	NE-CZ-NH1	7.57	124.08	120.30
2	K	13	DG	P-O3'-C3'	7.55	128.76	119.70
3	H	14	DC	O4'-C1'-N1	-7.55	102.72	108.00
1	F	170	ARG	NE-CZ-NH2	-7.37	116.61	120.30
3	L	2	DT	O4'-C1'-N1	-7.37	102.84	108.00
2	I	13	DG	P-O3'-C3'	7.37	128.54	119.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	G	2	DT	P-O3'-C3'	7.37	128.54	119.70
1	F	170	ARG	NE-CZ-NH1	7.30	123.95	120.30
2	G	4	DC	O4'-C1'-N1	-7.22	102.94	108.00
1	D	133	ARG	NE-CZ-NH1	7.16	123.88	120.30
1	D	170	ARG	NE-CZ-NH2	-7.09	116.75	120.30
1	F	133	ARG	CD-NE-CZ	6.98	133.37	123.60
3	J	2	DT	C5-C4-O4	-6.90	120.07	124.90
3	L	14	DC	O4'-C1'-N1	-6.86	103.19	108.00
2	G	13	DG	P-O3'-C3'	6.82	127.88	119.70
3	L	2	DT	N3-C4-O4	6.75	123.95	119.90
3	H	2	DT	O4'-C1'-N1	-6.72	103.30	108.00
2	I	4	DC	O4'-C1'-N1	-6.70	103.31	108.00
3	L	2	DT	C5-C4-O4	-6.58	120.30	124.90
3	J	2	DT	N3-C4-O4	6.52	123.81	119.90
3	L	6	DT	N3-C4-O4	6.51	123.81	119.90
3	J	14	DC	O4'-C1'-N1	-6.50	103.45	108.00
3	H	2	DT	C5-C4-O4	-6.50	120.35	124.90
3	H	13	DG	O4'-C1'-N9	-6.44	103.49	108.00
2	K	2	DT	P-O3'-C3'	6.42	127.41	119.70
1	D	154	ARG	NE-CZ-NH1	6.37	123.48	120.30
1	E	170	ARG	CD-NE-CZ	6.37	132.52	123.60
1	F	154	ARG	NE-CZ-NH1	6.36	123.48	120.30
2	I	2	DT	P-O3'-C3'	6.34	127.31	119.70
2	K	4	DC	O4'-C1'-N1	-6.33	103.57	108.00
3	H	2	DT	N3-C4-O4	6.30	123.68	119.90
3	H	10	DT	P-O3'-C3'	6.18	127.12	119.70
3	H	12	DT	C4-C5-C7	6.10	122.66	119.00
3	L	2	DT	C1'-O4'-C4'	-6.01	104.09	110.10
1	C	170	ARG	NE-CZ-NH2	-5.98	117.31	120.30
1	B	170	ARG	CG-CD-NE	5.79	123.96	111.80
3	L	10	DT	P-O3'-C3'	5.77	126.63	119.70
3	L	6	DT	C5-C4-O4	-5.74	120.89	124.90
1	C	154	ARG	NH1-CZ-NH2	-5.72	113.11	119.40
2	K	8	DA	O4'-C1'-N9	-5.72	104.00	108.00
2	K	12	DT	O4'-C1'-C2'	-5.71	101.33	105.90
3	L	13	DG	O4'-C1'-N9	-5.70	104.01	108.00
3	H	11	DA	P-O3'-C3'	5.70	126.54	119.70
1	A	170	ARG	NE-CZ-NH2	-5.68	117.46	120.30
3	J	2	DT	C1'-O4'-C4'	-5.54	104.56	110.10
2	G	8	DA	O4'-C1'-N9	-5.46	104.18	108.00
2	G	9	DC	O4'-C1'-N1	-5.41	104.21	108.00
3	H	2	DT	C1'-O4'-C4'	-5.28	104.82	110.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	J	13	DG	O4'-C1'-N9	-5.20	104.36	108.00
3	L	12	DT	C3'-C2'-C1'	-5.19	96.27	102.50
3	H	7	DC	O4'-C1'-N1	5.15	111.61	108.00
1	F	154	ARG	NH1-CZ-NH2	-5.14	113.74	119.40
2	K	9	DC	O4'-C1'-N1	-5.14	104.40	108.00
3	J	4	DC	N3-C2-O2	5.13	125.49	121.90
2	I	3	DG	P-O3'-C3'	5.04	125.75	119.70
2	G	3	DG	P-O3'-C3'	5.04	125.74	119.70
1	D	154	ARG	NH1-CZ-NH2	-5.01	113.89	119.40

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	1075	0	1098	21	0
1	B	1075	0	1098	18	0
1	C	1075	0	1098	20	0
1	D	1117	0	1138	34	0
1	E	1112	0	1133	33	0
1	F	1117	0	1138	32	0
2	G	326	0	182	2	0
2	I	326	0	182	1	0
2	K	326	0	182	1	0
3	H	324	0	184	5	0
3	J	324	0	184	4	0
3	L	324	0	184	3	0
4	A	12	0	12	1	0
4	B	12	0	12	3	0
4	C	12	0	12	0	0
4	D	12	0	12	1	0
4	E	12	0	12	1	0
4	F	12	0	12	1	0
5	A	64	0	0	5	0
5	B	59	0	0	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	C	66	0	0	4	0
5	D	47	0	0	4	0
5	E	44	0	0	4	0
5	F	45	0	0	1	0
5	G	18	0	0	0	0
5	H	25	0	0	0	0
5	I	13	0	0	0	0
5	J	20	0	0	0	0
5	K	19	0	0	0	0
5	L	15	0	0	0	0
All	All	9028	0	7873	157	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 10.

All (157) 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:F:87:VAL:HB	1:F:88:ARG:HD2	1.56	0.87
1:D:87:VAL:HB	1:D:88:ARG:HD2	1.56	0.85
1:E:87:VAL:HB	1:E:88:ARG:HD2	1.57	0.85
1:A:17:ASN:HB2	5:A:241:HOH:O	1.81	0.79
1:D:125:HIS:HD2	5:D:216:HOH:O	1.67	0.76
1:A:153:ALA:HB2	1:A:162:LEU:HD22	1.68	0.76
1:C:153:ALA:HB2	1:C:162:LEU:HD22	1.68	0.75
1:D:165:MET:HE2	5:D:240:HOH:O	1.86	0.74
1:B:153:ALA:HB2	1:B:162:LEU:HD22	1.69	0.73
1:E:165:MET:HE2	5:E:265:HOH:O	1.90	0.72
1:B:99:ARG:HD2	5:B:203:HOH:O	1.90	0.71
1:D:165:MET:HG3	5:D:240:HOH:O	1.89	0.71
1:C:99:ARG:HD2	5:C:216:HOH:O	1.91	0.69
1:E:24:ARG:O	1:E:28:ILE:HG13	1.93	0.68
1:A:99:ARG:HD2	5:A:298:HOH:O	1.94	0.67
1:E:132:ASP:OD1	4:E:200:ARG:HG3	1.94	0.67
1:D:18:ARG:HG3	1:D:21:ARG:NH1	2.10	0.67
1:F:24:ARG:O	1:F:28:ILE:HG13	1.95	0.66
1:E:165:MET:HG3	5:E:265:HOH:O	1.95	0.64
1:F:18:ARG:HG3	1:F:21:ARG:NH1	2.12	0.64
1:D:24:ARG:O	1:D:28:ILE:HG13	1.97	0.64
1:C:66:VAL:HG11	1:E:68:LEU:HD12	1.80	0.64
1:E:125:HIS:HD2	5:E:217:HOH:O	1.80	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:154:ARG:HH11	1:E:154:ARG:HG3	1.64	0.62
1:B:82:GLU:N	1:F:88:ARG:HB2	2.14	0.62
1:C:29:LEU:HD12	1:C:80:VAL:HG12	1.82	0.61
1:C:92:GLY:N	1:C:93:GLY:HA3	2.16	0.61
1:E:18:ARG:HG3	1:E:21:ARG:NH1	2.16	0.61
1:C:82:GLU:N	1:E:88:ARG:HB2	2.17	0.60
1:B:81:PRO:O	1:B:82:GLU:HB2	2.02	0.60
1:A:138:GLN:HG3	5:A:206:HOH:O	2.01	0.60
1:D:143:ILE:HG13	1:F:150:LEU:HD22	1.84	0.59
1:A:29:LEU:HD12	1:A:80:VAL:HG12	1.84	0.58
1:C:81:PRO:O	1:C:82:GLU:HB2	2.02	0.58
1:E:69:ARG:HD3	1:E:74:GLY:O	2.03	0.58
1:B:92:GLY:N	1:B:93:GLY:HA3	2.17	0.58
1:A:81:PRO:O	1:A:82:GLU:HB2	2.03	0.58
1:D:69:ARG:HD3	1:D:74:GLY:O	2.04	0.58
1:A:108:THR:HG22	1:A:117:LEU:HD22	1.86	0.57
1:D:105:LEU:HB3	1:D:170:ARG:HG3	1.87	0.57
1:B:29:LEU:HD12	1:B:80:VAL:HG12	1.85	0.57
1:A:92:GLY:N	1:A:93:GLY:HA3	2.18	0.57
1:B:108:THR:HG22	1:B:117:LEU:HD22	1.86	0.56
1:A:82:GLU:N	1:D:88:ARG:HB2	2.20	0.56
1:F:69:ARG:HD3	1:F:74:GLY:O	2.07	0.55
1:C:108:THR:HG22	1:C:117:LEU:HD22	1.88	0.55
1:D:142:THR:HG22	1:D:151:VAL:HG13	1.88	0.55
1:B:81:PRO:CB	1:F:88:ARG:HD3	2.36	0.55
1:B:22:GLN:HB3	1:B:63:LEU:HD11	1.89	0.55
1:B:81:PRO:HB2	1:F:88:ARG:HD3	1.88	0.55
1:F:133:ARG:O	1:F:133:ARG:HD3	2.06	0.54
1:D:129:SER:HB2	4:D:200:ARG:HD3	1.89	0.54
1:D:118:ARG:NH2	1:D:147:ASP:CG	2.61	0.54
3:J:2:DT:H2'	3:J:3:DG:C8	2.43	0.54
1:E:150:LEU:HD22	1:F:143:ILE:HG13	1.89	0.53
1:E:21:ARG:CD	1:E:51:VAL:HG11	2.38	0.53
1:F:21:ARG:CD	1:F:51:VAL:HG11	2.39	0.53
1:D:21:ARG:CD	1:D:51:VAL:HG11	2.39	0.52
1:E:82:GLU:C	1:E:84:GLY:H	2.12	0.52
1:B:82:GLU:H	1:F:88:ARG:HB2	1.74	0.52
1:F:105:LEU:HB3	1:F:170:ARG:HG3	1.90	0.52
3:H:2:DT:H2'	3:H:3:DG:C8	2.44	0.52
1:D:34:VAL:HG11	1:D:40:LEU:HB2	1.92	0.52
1:F:118:ARG:NH2	1:F:147:ASP:CG	2.63	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:21:ARG:HD3	1:F:51:VAL:HG11	1.92	0.51
3:L:2:DT:H2'	3:L:3:DG:C8	2.44	0.51
1:E:21:ARG:HD3	1:E:51:VAL:HG11	1.91	0.51
1:D:21:ARG:HD3	1:D:51:VAL:HG11	1.91	0.51
1:C:22:GLN:HB3	1:C:63:LEU:HD11	1.92	0.51
1:F:82:GLU:C	1:F:84:GLY:H	2.14	0.51
1:F:119:THR:HB	1:F:120:PRO:HD2	1.92	0.51
1:C:81:PRO:HB2	1:E:88:ARG:HD3	1.93	0.51
1:D:119:THR:HB	1:D:120:PRO:HD2	1.93	0.51
1:B:125:HIS:HB3	4:B:200:ARG:HH21	1.77	0.50
1:D:58:ARG:O	1:D:62:GLU:HG3	2.11	0.50
1:D:82:GLU:C	1:D:84:GLY:H	2.14	0.50
1:E:118:ARG:NH2	1:E:147:ASP:CG	2.65	0.50
1:F:132:ASP:OD1	4:F:200:ARG:N	2.44	0.50
1:D:118:ARG:HH22	1:D:147:ASP:CG	2.15	0.49
1:B:95:ASP:OD2	1:B:95:ASP:N	2.40	0.49
1:A:22:GLN:HB3	1:A:63:LEU:HD11	1.93	0.49
1:C:81:PRO:CB	1:E:88:ARG:HD3	2.42	0.49
5:B:216:HOH:O	1:C:155:GLU:HG3	2.13	0.49
1:E:142:THR:HG22	1:E:151:VAL:HG13	1.95	0.49
1:F:58:ARG:O	1:F:62:GLU:HG3	2.14	0.48
1:A:81:PRO:HB2	1:D:88:ARG:HD3	1.95	0.48
1:E:58:ARG:O	1:E:62:GLU:HG3	2.13	0.48
3:J:2:DT:H2''	3:J:3:DG:H5'	1.95	0.48
1:A:148:THR:HG1	4:B:200:ARG:N	2.12	0.48
1:C:80:VAL:HA	1:C:81:PRO:HD3	1.66	0.47
1:E:34:VAL:HG11	1:E:40:LEU:HB2	1.97	0.47
1:A:35:ARG:HG3	5:A:294:HOH:O	2.14	0.47
1:F:34:VAL:HG11	1:F:40:LEU:HB2	1.97	0.47
1:E:119:THR:HB	1:E:120:PRO:HD2	1.96	0.47
1:F:118:ARG:HH22	1:F:147:ASP:CG	2.18	0.47
1:C:35:ARG:HG3	5:C:215:HOH:O	2.15	0.46
1:C:95:ASP:OD2	1:C:95:ASP:N	2.45	0.46
1:D:68:LEU:HD12	1:D:79:VAL:HG21	1.96	0.46
3:J:2:DT:H2''	3:J:3:DG:C5'	2.44	0.46
1:E:118:ARG:HH22	1:E:147:ASP:CG	2.19	0.46
1:C:33:GLN:HE21	1:E:133:ARG:HH11	1.64	0.46
1:B:125:HIS:HB3	4:B:200:ARG:NH2	2.31	0.45
1:F:90:VAL:C	1:F:92:GLY:H	2.19	0.45
1:F:142:THR:HG22	1:F:151:VAL:HG13	1.98	0.45
1:A:80:VAL:HA	1:A:81:PRO:HD3	1.65	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:35:ARG:HH12	1:F:75:THR:HB	1.82	0.45
3:H:1:DT:HO5'	3:H:1:DT:H6	1.64	0.45
3:H:2:DT:H2''	3:H:3:DG:C5'	2.46	0.45
1:A:33:GLN:HE21	1:D:133:ARG:HH11	1.64	0.44
3:L:2:DT:H2''	3:L:3:DG:C5'	2.48	0.44
1:F:119:THR:HB	1:F:120:PRO:CD	2.47	0.44
1:F:35:ARG:NH1	1:F:75:THR:HB	2.33	0.44
1:E:150:LEU:HD12	1:E:150:LEU:HA	1.82	0.44
1:A:81:PRO:CB	1:D:88:ARG:HD3	2.47	0.44
1:C:82:GLU:H	1:E:88:ARG:HB2	1.81	0.44
1:E:119:THR:HB	1:E:120:PRO:CD	2.47	0.44
1:E:153:ALA:HB2	1:E:162:LEU:HD22	2.00	0.44
1:B:125:HIS:HD2	5:B:222:HOH:O	2.01	0.43
1:D:68:LEU:HD23	1:D:68:LEU:HA	1.74	0.43
1:E:24:ARG:HG3	1:E:49:ILE:HD11	2.00	0.43
3:J:1:DT:H6	3:J:1:DT:HO5'	1.64	0.43
1:B:138:GLN:HG3	5:B:207:HOH:O	2.18	0.43
1:A:100:LEU:HD23	1:A:100:LEU:HA	1.80	0.43
1:E:90:VAL:C	1:E:92:GLY:H	2.21	0.43
1:D:133:ARG:HD2	1:D:133:ARG:HA	1.69	0.43
1:D:24:ARG:HG3	1:D:49:ILE:HD11	2.00	0.42
2:G:4:DC:H2''	2:G:5:DA:C8	2.54	0.42
3:L:7:DC:H2''	3:L:8:DG:C8	2.54	0.42
1:F:69:ARG:HA	5:F:339:HOH:O	2.18	0.42
1:F:150:LEU:HD12	1:F:150:LEU:HA	1.81	0.42
1:B:38:ASN:HB3	5:E:218:HOH:O	2.20	0.42
1:D:119:THR:HB	1:D:120:PRO:CD	2.49	0.42
1:F:24:ARG:HG3	1:F:49:ILE:HD11	1.99	0.42
1:A:155:GLU:HG3	5:C:212:HOH:O	2.19	0.42
1:D:125:HIS:CD2	5:D:216:HOH:O	2.53	0.42
1:A:98:ALA:HA	1:A:101:LEU:HD12	2.02	0.42
1:D:90:VAL:C	1:D:92:GLY:H	2.22	0.42
1:D:77:ILE:O	1:D:77:ILE:HG13	2.20	0.42
3:H:2:DT:H2''	3:H:3:DG:H5'	2.02	0.41
1:C:62:GLU:OE1	5:C:217:HOH:O	2.20	0.41
1:C:150:LEU:HA	1:C:150:LEU:HD12	1.81	0.41
3:H:2:DT:H2''	3:H:3:DG:O5'	2.20	0.41
1:D:35:ARG:HH12	1:D:75:THR:HB	1.86	0.41
1:F:143:ILE:HD12	1:F:143:ILE:HG23	1.71	0.41
1:B:98:ALA:HA	1:B:101:LEU:HD12	2.03	0.41
1:D:150:LEU:HD22	1:E:143:ILE:HG13	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:K:6:DT:H2''	2:K:7:DA:C8	2.56	0.41
1:C:100:LEU:HD23	1:C:100:LEU:HA	1.88	0.41
1:F:77:ILE:O	1:F:77:ILE:HG13	2.21	0.41
1:E:35:ARG:HH12	1:E:75:THR:HB	1.86	0.41
1:F:68:LEU:HD12	1:F:79:VAL:HG21	2.02	0.41
1:A:17:ASN:CB	5:A:241:HOH:O	2.52	0.40
1:A:129:SER:CB	4:A:200:ARG:HD3	2.51	0.40
1:D:24:ARG:HH11	1:D:24:ARG:HD2	1.78	0.40
2:I:4:DC:H2''	2:I:5:DA:C8	2.56	0.40
2:G:6:DT:H2''	2:G:7:DA:C8	2.57	0.40
1:E:35:ARG:NH1	1:E:75:THR:HB	2.36	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	145/170 (85%)	144 (99%)	1 (1%)	0	100	100
1	B	145/170 (85%)	144 (99%)	1 (1%)	0	100	100
1	C	145/170 (85%)	143 (99%)	2 (1%)	0	100	100
1	D	153/170 (90%)	143 (94%)	8 (5%)	2 (1%)	12	12
1	E	152/170 (89%)	142 (93%)	8 (5%)	2 (1%)	12	12
1	F	153/170 (90%)	143 (94%)	8 (5%)	2 (1%)	12	12
All	All	893/1020 (88%)	859 (96%)	28 (3%)	6 (1%)	22	26

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	168	ASN

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Mol	Chain	Res	Type
1	F	168	ASN
1	D	87	VAL
1	E	87	VAL
1	E	168	ASN
1	F	87	VAL

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	108/122 (88%)	100 (93%)	8 (7%)	13	17
1	B	108/122 (88%)	100 (93%)	8 (7%)	13	17
1	C	108/122 (88%)	100 (93%)	8 (7%)	13	17
1	D	113/122 (93%)	103 (91%)	10 (9%)	10	12
1	E	113/122 (93%)	102 (90%)	11 (10%)	8	9
1	F	113/122 (93%)	103 (91%)	10 (9%)	10	12
All	All	663/732 (91%)	608 (92%)	55 (8%)	11	14

All (55) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	63	LEU
1	A	68	LEU
1	A	75	THR
1	A	80	VAL
1	A	95	ASP
1	A	105	LEU
1	A	150	LEU
1	A	162	LEU
1	B	63	LEU
1	B	68	LEU
1	B	75	THR
1	B	80	VAL
1	B	95	ASP

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Mol	Chain	Res	Type
1	B	105	LEU
1	B	150	LEU
1	B	162	LEU
1	C	63	LEU
1	C	68	LEU
1	C	75	THR
1	C	80	VAL
1	C	95	ASP
1	C	105	LEU
1	C	150	LEU
1	C	162	LEU
1	D	63	LEU
1	D	72	ASP
1	D	79	VAL
1	D	80	VAL
1	D	85	SER
1	D	87	VAL
1	D	88	ARG
1	D	94	THR
1	D	105	LEU
1	D	138	GLN
1	E	63	LEU
1	E	68	LEU
1	E	72	ASP
1	E	79	VAL
1	E	80	VAL
1	E	85	SER
1	E	87	VAL
1	E	88	ARG
1	E	94	THR
1	E	105	LEU
1	E	138	GLN
1	F	63	LEU
1	F	79	VAL
1	F	80	VAL
1	F	85	SER
1	F	87	VAL
1	F	88	ARG
1	F	94	THR
1	F	105	LEU
1	F	133	ARG
1	F	138	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	33	GLN
1	D	168	ASN
1	E	168	ASN
1	F	168	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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

6 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	ARG	E	200	-	10,11,11	0.68	0	11,13,13	1.16	1 (9%)
4	ARG	A	200	-	10,11,11	0.79	0	11,13,13	0.78	0
4	ARG	C	200	-	10,11,11	0.66	0	11,13,13	0.73	0
4	ARG	F	200	-	10,11,11	0.83	0	11,13,13	0.72	0
4	ARG	B	200	-	10,11,11	0.63	0	11,13,13	0.98	1 (9%)
4	ARG	D	200	-	10,11,11	0.73	0	11,13,13	1.06	1 (9%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	ARG	E	200	-	-	2/11/11/11	-
4	ARG	A	200	-	-	1/11/11/11	-
4	ARG	C	200	-	-	0/11/11/11	-
4	ARG	F	200	-	-	0/11/11/11	-
4	ARG	B	200	-	-	3/11/11/11	-
4	ARG	D	200	-	-	0/11/11/11	-

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	E	200	ARG	OXT-C-CA	2.54	122.04	113.38
4	D	200	ARG	OXT-C-CA	2.37	121.47	113.38
4	B	200	ARG	OXT-C-CA	2.16	120.75	113.38

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	200	ARG	N-CA-CB-CG
4	B	200	ARG	C-CA-CB-CG
4	B	200	ARG	NE-CD-CG-CB
4	E	200	ARG	CA-CB-CG-CD
4	A	200	ARG	OXT-C-CA-N
4	E	200	ARG	N-CA-CB-CG

There are no ring outliers.

5 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	E	200	ARG	1	0
4	A	200	ARG	1	0
4	F	200	ARG	1	0
4	B	200	ARG	3	0
4	D	200	ARG	1	0

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	A	149/170 (87%)	0.10	4 (2%) 54 62	32, 47, 81, 114	0
1	B	149/170 (87%)	-0.00	2 (1%) 77 81	32, 47, 81, 113	0
1	C	149/170 (87%)	0.03	2 (1%) 77 81	31, 46, 81, 115	0
1	D	155/170 (91%)	0.53	15 (9%) 7 10	31, 57, 132, 165	0
1	E	154/170 (90%)	0.48	13 (8%) 11 15	32, 58, 131, 164	0
1	F	155/170 (91%)	0.46	11 (7%) 16 21	32, 57, 132, 165	0
2	G	16/16 (100%)	-0.68	0 100 100	43, 61, 73, 74	0
2	I	16/16 (100%)	-0.63	0 100 100	43, 59, 73, 75	0
2	K	16/16 (100%)	-0.66	0 100 100	43, 60, 74, 75	0
3	H	16/16 (100%)	-0.66	0 100 100	41, 63, 71, 73	0
3	J	16/16 (100%)	-0.70	0 100 100	41, 63, 71, 73	0
3	L	16/16 (100%)	-0.74	0 100 100	42, 63, 72, 74	0
All	All	1007/1116 (90%)	0.18	47 (4%) 31 38	31, 52, 99, 165	0

All (47) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	85	SER	10.3
1	E	85	SER	8.7
1	E	87	VAL	8.7
1	E	84	GLY	8.4
1	D	88	ARG	7.7
1	F	85	SER	7.6
1	A	90	VAL	7.5
1	F	84	GLY	6.9
1	D	90	VAL	6.7
1	D	87	VAL	6.5
1	B	90	VAL	6.4

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Mol	Chain	Res	Type	RSRZ
1	F	88	ARG	6.3
1	E	88	ARG	6.0
1	E	91	SER	5.8
1	C	90	VAL	5.6
1	F	87	VAL	5.5
1	F	91	SER	5.3
1	F	83	ASP	5.2
1	D	91	SER	5.1
1	D	83	ASP	4.9
1	D	84	GLY	4.6
1	E	83	ASP	4.5
1	F	82	GLU	4.2
1	B	15	ALA	3.7
1	C	15	ALA	3.3
1	F	90	VAL	3.3
1	D	86	PRO	3.3
1	E	90	VAL	3.3
1	E	72	ASP	3.1
1	A	15	ALA	2.9
1	A	92	GLY	2.8
1	E	73	GLY	2.8
1	D	82	GLU	2.6
1	E	82	GLU	2.6
1	D	89	GLY	2.6
1	F	72	ASP	2.5
1	D	92	GLY	2.4
1	D	50	GLU	2.4
1	D	72	ASP	2.3
1	D	71	ALA	2.3
1	F	43	LEU	2.2
1	A	16	ALA	2.1
1	D	48	GLY	2.1
1	E	74	GLY	2.1
1	F	92	GLY	2.1
1	E	86	PRO	2.1
1	E	46	ALA	2.1

6.2 Non-standard residues in protein, DNA, RNA chains

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

6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

6.4 Ligands [i](#)

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(Å ²)	Q<0.9
4	ARG	D	200	12/12	0.81	0.45	47,59,73,82	12
4	ARG	F	200	12/12	0.85	0.34	54,61,72,81	12
4	ARG	E	200	12/12	0.86	0.33	45,58,76,84	12
4	ARG	B	200	12/12	0.89	0.31	40,54,73,82	12
4	ARG	A	200	12/12	0.89	0.44	38,56,76,84	12
4	ARG	C	200	12/12	0.91	0.32	37,57,76,77	12

6.5 Other polymers [i](#)

There are no such residues in this entry.