



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 04:51 AM UTC

PDB ID : 4LLL / pdb_00004lll
Title : Crystal structure of S. aureus MepR-DNA complex
Authors : Birukou, I.; Brennan, R.G.
Deposited on : 2013-07-09
Resolution : 3.04 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:

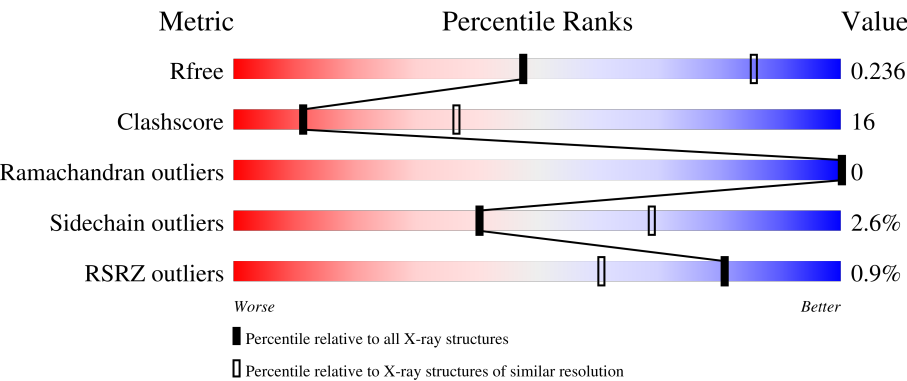
MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
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 3.04 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	3685 (3.08-3.00)
Clashscore	190562	4007 (3.08-3.00)
Ramachandran outliers	187476	3834 (3.08-3.00)
Sidechain outliers	187428	3836 (3.08-3.00)
RSRZ outliers	180081	3684 (3.08-3.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments 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	140	55%41%..
1	B	140	% 69%30%.
1	C	140	60%35%..
1	D	140	% 67%29%..
1	I	140	% 76%21%..

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Mol	Chain	Length	Quality of chain
1	J	140	2%64%31%. .
1	M	140	%64%24%11%
1	O	140	%69%27%. .
2	E	24	50%50%
2	F	24	54%38%8%
2	G	24	33%58%8%
2	H	24	42%58%
2	K	24	58%38%.
2	L	24	46%54%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 11037 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 MepR.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	138	Total	C	N	O	S	0	0	0
			1080	675	190	210	5			
1	B	138	Total	C	N	O	S	0	0	0
			1091	683	191	213	4			
1	C	138	Total	C	N	O	S	0	0	0
			1051	659	181	206	5			
1	D	137	Total	C	N	O	S	0	0	0
			1041	656	183	197	5			
1	I	137	Total	C	N	O	S	0	0	0
			979	610	169	197	3			
1	J	138	Total	C	N	O	S	0	0	0
			1008	630	173	201	4			
1	M	124	Total	C	N	O	S	0	0	0
			900	558	163	174	5			
1	O	135	Total	C	N	O	S	0	0	0
			953	594	171	184	4			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	SER	-	expression tag	UNP Q5Y812
A	1	ASN	-	expression tag	UNP Q5Y812
B	0	SER	-	expression tag	UNP Q5Y812
B	1	ASN	-	expression tag	UNP Q5Y812
C	0	SER	-	expression tag	UNP Q5Y812
C	1	ASN	-	expression tag	UNP Q5Y812
D	0	SER	-	expression tag	UNP Q5Y812
D	1	ASN	-	expression tag	UNP Q5Y812
I	0	SER	-	expression tag	UNP Q5Y812
I	1	ASN	-	expression tag	UNP Q5Y812
J	0	SER	-	expression tag	UNP Q5Y812
J	1	ASN	-	expression tag	UNP Q5Y812
M	0	SER	-	expression tag	UNP Q5Y812

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Chain	Residue	Modelled	Actual	Comment	Reference
M	1	ASN	-	expression tag	UNP Q5Y812
O	0	SER	-	expression tag	UNP Q5Y812
O	1	ASN	-	expression tag	UNP Q5Y812

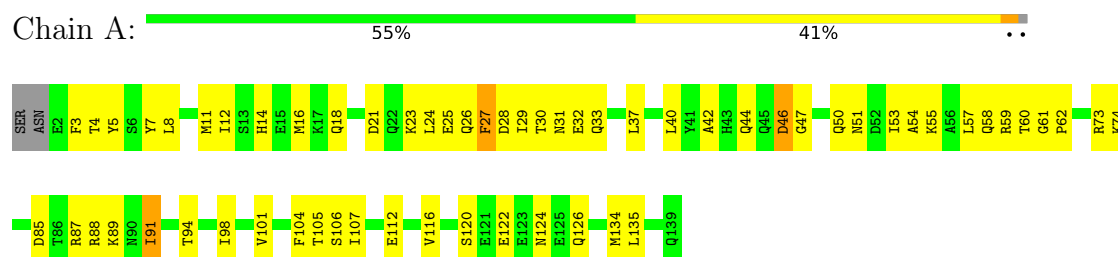
- Molecule 2 is a DNA chain called Palindromized mepR operator sequence.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	24	Total 489	C 238	N 86	O 142	P 23	0	0	0
2	F	24	Total 489	C 238	N 86	O 142	P 23	0	0	0
2	G	24	Total 489	C 238	N 86	O 142	P 23	0	0	0
2	H	24	Total 489	C 238	N 86	O 142	P 23	0	0	0
2	K	24	Total 489	C 238	N 86	O 142	P 23	0	0	0
2	L	24	Total 489	C 238	N 86	O 142	P 23	0	0	0

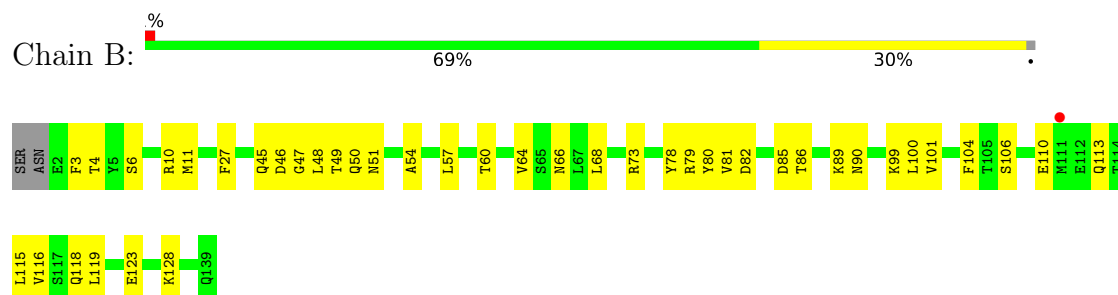
3 Residue-property plots [i](#)

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.

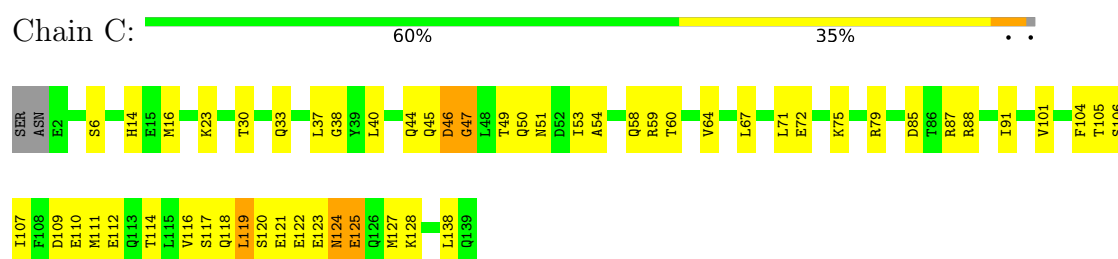
• Molecule 1: MepR



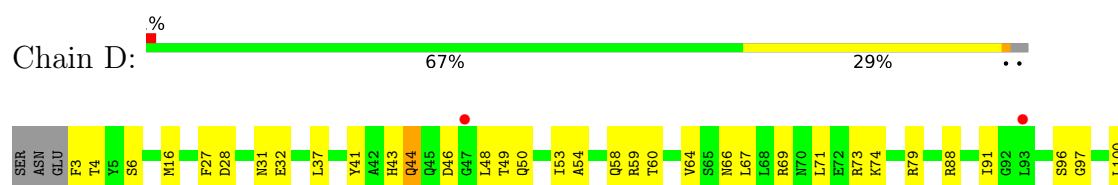
• Molecule 1: MepR

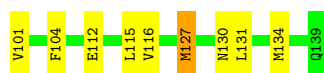


• Molecule 1: MepR

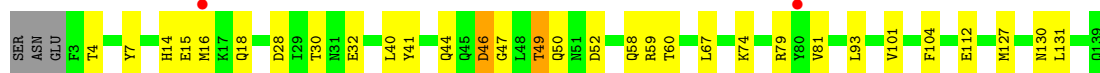
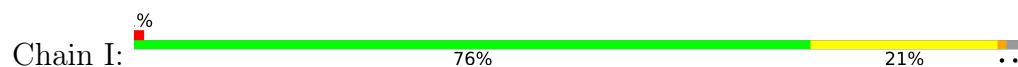


• Molecule 1: MepR

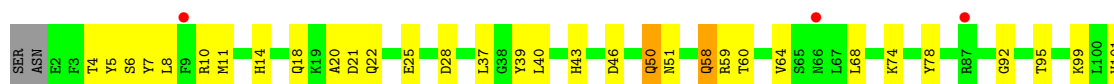




- Molecule 1: MepR



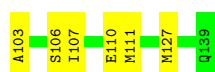
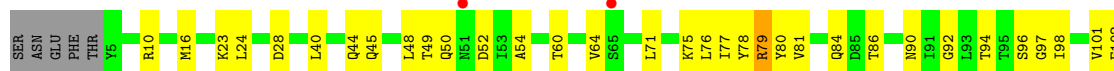
- Molecule 1: MepR



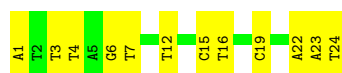
- Molecule 1: MepR



- Molecule 1: MepR

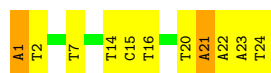


- Molecule 2: Palindromized mepR operator sequence

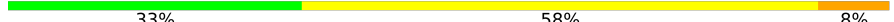


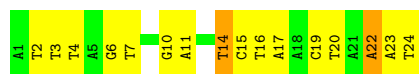
- Molecule 2: Palindromized mepR operator sequence

Chain F:  54% 38% 8%



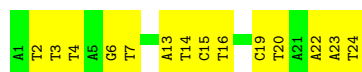
- Molecule 2: Palindromized mepR operator sequence

Chain G:  33% 58% 8%



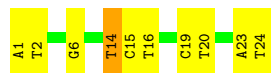
- Molecule 2: Palindromized mepR operator sequence

Chain H:  42% 58%

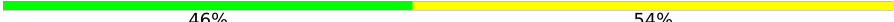


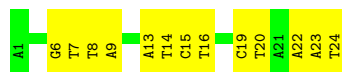
- Molecule 2: Palindromized mepR operator sequence

Chain K:  58% 38% 4%



- Molecule 2: Palindromized mepR operator sequence

Chain L:  46% 54%



4 Data and refinement statistics

Property	Value	Source
Space group	P 31	Depositor
Cell constants a, b, c, α , β , γ	130.18Å 130.18Å 124.71Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	42.61 – 3.04 42.61 – 3.04	Depositor EDS
% Data completeness (in resolution range)	99.6 (42.61-3.04) 96.4 (42.61-3.04)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.12 (at 3.06Å)	Xtriage
Refinement program	PHENIX 1.8.2_1309	Depositor
R, R_{free}	0.194 , 0.234 0.195 , 0.236	Depositor DCC
R_{free} test set	1931 reflections (4.24%)	wwPDB-VP
Wilson B-factor (Å ²)	76.4	Xtriage
Anisotropy	0.149	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.25 , 41.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.041 for -h,-k,l 0.067 for h,-h-k,-l 0.047 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	11037	wwPDB-VP
Average B, all atoms (Å ²)	75.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 21.41 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 7.0597e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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	0.87	2/1094 (0.2%)	1.20	13/1477 (0.9%)
1	B	0.58	0/1105	0.99	3/1489 (0.2%)
1	C	0.70	2/1065 (0.2%)	1.28	17/1443 (1.2%)
1	D	0.62	0/1055	0.95	1/1425 (0.1%)
1	I	0.46	0/992	0.95	1/1354 (0.1%)
1	J	0.45	0/1021	1.06	6/1387 (0.4%)
1	M	0.47	0/910	0.88	2/1234 (0.2%)
1	O	0.47	0/964	0.93	2/1309 (0.2%)
2	E	0.49	0/548	1.19	3/844 (0.4%)
2	F	0.49	0/548	1.23	4/844 (0.5%)
2	G	0.56	0/548	1.31	3/844 (0.4%)
2	H	0.51	0/548	1.26	2/844 (0.2%)
2	K	0.31	0/548	0.96	1/844 (0.1%)
2	L	0.36	0/548	1.06	0/844
All	All	0.57	4/11494 (0.0%)	1.09	58/16182 (0.4%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	26	GLN	CA-C	7.19	1.62	1.52
1	C	123	GLU	CA-C	6.01	1.61	1.52
1	C	118	GLN	CA-C	5.39	1.60	1.52
1	A	62	PRO	N-CD	5.02	1.54	1.47

All (58) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	119	LEU	N-CA-C	16.03	134.66	108.52
1	A	27	PHE	N-CA-C	15.90	132.07	113.15
1	C	123	GLU	N-CA-C	10.95	126.47	112.89
1	A	25	GLU	N-CA-C	-10.86	99.22	111.71
1	C	117	SER	N-CA-C	10.66	124.19	111.82
1	J	138	LEU	N-CA-C	9.40	121.60	111.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	J	136	SER	N-CA-C	-9.27	101.17	111.28
1	C	124	ASN	N-CA-CB	9.21	123.92	110.20
1	A	61	GLY	O-C-N	8.81	124.24	121.07
1	I	46	ASP	N-CA-C	8.18	123.03	112.89
1	C	128	LYS	N-CA-C	-8.10	102.53	111.36
1	C	125	GLU	N-CA-C	-8.09	102.46	111.28
1	A	27	PHE	N-CA-CB	-7.96	98.21	110.46
1	C	120	SER	N-CA-CB	7.86	121.48	110.17
1	J	125	GLU	N-CA-C	-7.84	102.74	111.28
2	F	1	DA	C2'-C3'-O3'	-7.76	99.86	111.50
1	D	44	GLN	N-CA-C	7.60	120.53	111.33
1	A	46	ASP	N-CA-C	7.57	120.60	111.82
1	C	119	LEU	CB-CA-C	-7.40	98.05	110.19
1	J	122	GLU	N-CA-C	-6.91	104.97	113.19
1	C	120	SER	N-CA-C	-6.79	100.93	110.50
2	G	22	DA	O4'-C1'-N9	6.72	118.48	108.40
1	J	112	GLU	N-CA-C	-6.31	104.03	111.03
2	E	7	DT	C4'-C3'-O3'	-6.19	100.71	110.00
1	A	26	GLN	CA-C-O	6.18	128.25	121.02
1	O	10	ARG	N-CA-C	-6.10	104.19	111.69
1	C	118	GLN	CA-C-O	6.01	126.81	119.11
1	A	106	SER	N-CA-C	5.99	118.58	111.33
2	G	14	DT	O4'-C4'-C3'	-5.86	96.61	105.40
1	M	14	HIS	N-CA-C	5.82	117.30	111.07
1	A	105	THR	N-CA-C	5.73	117.53	111.28
1	A	16	MET	N-CA-C	-5.72	105.04	111.28
1	C	122	GLU	N-CA-C	5.70	117.17	111.07
1	A	91	ILE	CB-CA-C	-5.58	102.52	110.83
1	B	3	PHE	N-CA-C	5.55	117.69	109.69
1	A	42	ALA	N-CA-C	5.53	117.11	111.14
2	F	21	DA	N9-C1'-C2'	5.48	121.72	113.50
2	E	19	DC	O4'-C4'-C3'	-5.47	97.19	105.40
2	H	7	DT	C4'-C3'-O3'	-5.45	101.82	110.00
1	C	123	GLU	CB-CA-C	-5.45	98.60	109.99
1	C	46	ASP	N-CA-C	5.44	118.14	111.82
1	C	38	GLY	N-CA-C	-5.39	106.26	112.73
2	K	14	DT	O4'-C4'-C3'	-5.39	97.32	105.40
2	E	12	DT	O4'-C1'-C2'	-5.38	98.34	106.40
1	C	123	GLU	CA-C-O	5.31	125.90	119.31
2	H	4	DT	C4'-C3'-O3'	-5.29	102.06	110.00
1	M	46	ASP	N-CA-C	5.28	116.84	111.14
2	G	17	DA	O5'-C5'-C4'	-5.21	102.98	110.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	J	50	GLN	N-CA-C	5.19	116.93	111.28
2	F	21	DA	C4'-C3'-O3'	-5.15	102.27	110.00
1	C	47	GLY	N-CA-C	-5.15	98.37	113.30
1	B	66	ASN	N-CA-C	5.13	116.88	111.28
2	F	21	DA	C1'-O4'-C4'	-5.13	102.01	109.70
1	A	61	GLY	CA-C-N	-5.11	113.37	119.05
1	A	61	GLY	C-N-CA	-5.11	113.37	119.05
1	C	118	GLN	CB-CA-C	-5.05	99.90	109.95
1	B	73	ARG	N-CA-C	-5.04	105.98	111.82
1	O	84	GLN	N-CA-C	-5.02	107.11	113.43

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	1080	0	1035	39	0
1	B	1091	0	1057	38	0
1	C	1051	0	980	38	1
1	D	1041	0	984	40	1
1	I	979	0	841	31	0
1	J	1008	0	896	45	0
1	M	900	0	792	28	0
1	O	953	0	825	26	1
2	E	489	0	276	10	1
2	F	489	0	276	16	0
2	G	489	0	276	16	0
2	H	489	0	276	13	0
2	K	489	0	276	11	0
2	L	489	0	276	15	0
All	All	11037	0	9066	307	2

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 16.

All (307) close contacts within the same asymmetric unit are listed below, sorted by their clash

magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:1:DA:H2'	2:F:2:DT:H71	1.35	1.05
1:O:76:LEU:O	1:O:94:THR:HG22	1.58	1.03
1:I:127:MET:HG3	1:J:127:MET:HE2	1.42	0.99
1:A:28:ASP:O	1:A:74:LYS:NZ	1.96	0.99
1:D:127:MET:O	1:D:131:LEU:HD13	1.64	0.97
1:O:98:ILE:O	1:O:102:GLU:HG3	1.64	0.96
1:A:5:TYR:CD1	1:A:134:MET:HE1	2.04	0.93
1:B:85:ASP:HA	2:G:23:DA:H5'	1.51	0.92
1:O:97:GLY:O	1:O:101:VAL:HG23	1.72	0.90
1:A:5:TYR:CE1	1:A:134:MET:HE1	2.07	0.89
1:J:50:GLN:OE1	2:L:6:DG:H3'	1.73	0.88
1:I:58:GLN:O	1:I:59:ARG:NH1	2.07	0.88
1:M:49:THR:HG22	1:M:51:ASN:H	1.39	0.86
1:I:32:GLU:OE2	1:I:32:GLU:N	2.09	0.84
1:C:119:LEU:HD23	1:C:124:ASN:HA	1.60	0.83
1:O:76:LEU:C	1:O:94:THR:HG22	2.03	0.82
1:M:49:THR:HG22	1:M:50:GLN:N	1.94	0.82
1:J:4:THR:HG23	1:J:7:TYR:H	1.45	0.81
1:A:8:LEU:O	1:A:12:ILE:HG13	1.82	0.80
1:C:127:MET:HE3	1:D:131:LEU:CD1	2.12	0.80
1:B:50:GLN:HE22	2:H:6:DG:H3'	1.47	0.79
1:C:127:MET:HG3	1:D:127:MET:HE2	1.65	0.78
1:J:50:GLN:HG2	1:J:68:LEU:HD12	1.65	0.77
1:D:127:MET:HG3	1:D:131:LEU:HD13	1.64	0.77
1:I:127:MET:SD	1:J:127:MET:HG3	2.23	0.77
1:B:54:ALA:HA	1:B:64:VAL:HG21	1.68	0.76
1:C:127:MET:HE3	1:D:131:LEU:HD11	1.67	0.76
1:D:27:PHE:HB3	1:D:100:LEU:HD11	1.68	0.75
1:A:11:MET:HE1	1:B:57:LEU:O	1.85	0.75
1:J:50:GLN:HG2	1:J:68:LEU:CD1	2.21	0.71
1:J:124:ASN:OD1	1:J:125:GLU:N	2.24	0.71
1:B:4:THR:HG22	1:B:6:SER:H	1.55	0.70
1:J:121:GLU:C	1:J:123:GLU:H	1.97	0.70
1:A:4:THR:HG23	1:A:7:TYR:H	1.57	0.70
1:A:87:ARG:NH1	2:G:4:DT:H1'	2.07	0.69
1:B:49:THR:HG22	1:B:90:ASN:OD1	1.90	0.69
2:F:1:DA:C2'	2:F:2:DT:H71	2.19	0.69
1:C:127:MET:HG3	1:D:127:MET:CE	2.22	0.69
1:D:127:MET:HG3	1:D:131:LEU:CD1	2.21	0.69
1:C:6:SER:HA	1:D:16:MET:HE1	1.75	0.69
1:A:40:LEU:O	1:A:44:GLN:HB2	1.94	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:O:75:LYS:O	1:O:94:THR:HG21	1.94	0.67
1:I:28:ASP:O	1:I:74:LYS:NZ	2.26	0.67
1:B:81:VAL:O	1:B:81:VAL:HG13	1.94	0.67
1:M:60:THR:HG22	1:M:62:PRO:HD2	1.75	0.66
2:K:1:DA:H2'	2:K:2:DT:C6	2.31	0.66
1:I:15:GLU:OE2	1:J:58:GLN:NE2	2.29	0.65
1:J:39:TYR:HE1	1:J:43:HIS:CE1	2.14	0.65
1:D:69:ARG:HD3	1:D:73:ARG:CZ	2.27	0.65
1:J:39:TYR:CE1	1:J:43:HIS:CE1	2.85	0.65
1:B:119:LEU:HD23	1:B:123:GLU:HB3	1.79	0.64
1:M:54:ALA:HA	1:M:64:VAL:HG21	1.78	0.64
1:O:40:LEU:O	1:O:44:GLN:HB2	1.96	0.64
1:I:50:GLN:OE1	1:I:79:ARG:NH1	2.28	0.64
1:B:50:GLN:NE2	2:H:6:DG:H3'	2.12	0.64
1:O:79:ARG:HA	1:O:90:ASN:O	1.98	0.64
1:D:43:HIS:HB2	1:D:48:LEU:HD11	1.79	0.64
1:C:23:LYS:HB3	1:C:107:ILE:HD13	1.80	0.63
2:L:19:DC:H2''	2:L:20:DT:H5'	1.80	0.63
1:D:127:MET:O	1:D:131:LEU:CD1	2.43	0.63
1:A:3:PHE:HB2	1:A:8:LEU:HD21	1.80	0.63
1:A:134:MET:HE2	1:B:115:LEU:HD22	1.80	0.63
1:J:121:GLU:C	1:J:123:GLU:N	2.57	0.63
2:K:1:DA:H2''	2:K:2:DT:O5'	1.99	0.62
1:A:53:ILE:HG22	1:A:57:LEU:HD12	1.82	0.61
1:M:49:THR:CG2	1:M:50:GLN:N	2.63	0.61
1:A:89:LYS:NZ	2:G:6:DG:O3'	2.34	0.61
1:I:127:MET:CG	1:J:127:MET:HE2	2.25	0.61
1:M:49:THR:HG22	1:M:50:GLN:H	1.65	0.61
2:L:7:DT:H2'	2:L:8:DT:H71	1.83	0.61
1:A:50:GLN:HG2	1:A:91:ILE:HD11	1.84	0.60
1:J:20:ALA:HA	1:J:111:MET:HE1	1.83	0.60
1:J:59:ARG:NH1	2:K:14:DT:OP1	2.34	0.60
1:C:110:GLU:O	1:C:114:THR:OG1	2.18	0.60
1:A:32:GLU:N	1:A:32:GLU:OE1	2.32	0.59
1:M:49:THR:HG22	1:M:51:ASN:N	2.16	0.59
1:M:40:LEU:HB3	1:M:93:LEU:HD13	1.84	0.59
1:J:20:ALA:CA	1:J:111:MET:HE1	2.32	0.59
1:B:106:SER:O	1:B:110:GLU:HG2	2.03	0.58
1:A:23:LYS:HB3	1:A:107:ILE:HD13	1.86	0.58
2:L:8:DT:H2''	2:L:9:DA:OP2	2.03	0.58
1:A:14:HIS:O	1:A:18:GLN:HG2	2.03	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:50:GLN:HG2	1:B:68:LEU:CD1	2.33	0.58
1:C:14:HIS:NE2	2:E:15:DC:OP1	2.33	0.58
1:J:28:ASP:OD2	1:J:74:LYS:HE3	2.04	0.58
1:I:127:MET:HG3	1:J:127:MET:CE	2.27	0.57
1:O:16:MET:HA	1:O:111:MET:HE1	1.85	0.57
1:A:30:THR:OG1	2:H:13:DA:OP1	2.17	0.57
1:C:60:THR:HG23	2:F:15:DC:OP2	2.04	0.57
1:O:78:TYR:O	1:O:92:GLY:N	2.29	0.57
1:I:127:MET:HA	1:J:127:MET:HE2	1.86	0.57
2:L:9:DA:OP2	2:L:9:DA:H2'	2.05	0.57
2:H:22:DA:H2''	2:H:23:DA:C8	2.40	0.56
2:E:15:DC:H2'	2:E:16:DT:C6	2.41	0.56
1:B:78:TYR:HE1	1:B:80:TYR:HD2	1.54	0.56
2:K:23:DA:H2''	2:K:24:DT:H5'	1.88	0.56
2:E:1:DA:H61	2:F:24:DT:H3	1.52	0.56
1:C:121:GLU:O	1:C:125:GLU:CB	2.54	0.56
1:D:60:THR:HG23	2:E:15:DC:P	2.47	0.55
1:J:135:LEU:O	1:J:136:SER:C	2.47	0.55
1:B:85:ASP:HA	2:G:23:DA:C5'	2.32	0.55
1:B:113:GLN:HA	1:B:116:VAL:HG22	1.89	0.55
1:C:40:LEU:O	1:C:44:GLN:HB2	2.07	0.54
1:O:98:ILE:O	1:O:102:GLU:CG	2.49	0.54
1:M:127:MET:SD	1:O:127:MET:HA	2.47	0.54
1:A:51:ASN:O	1:A:55:LYS:HG3	2.07	0.54
2:L:19:DC:H2'	2:L:20:DT:H71	1.90	0.54
1:C:104:PHE:CD1	1:C:104:PHE:C	2.85	0.54
1:A:94:THR:O	1:A:98:ILE:HG13	2.08	0.54
1:I:40:LEU:O	1:I:44:GLN:HB2	2.08	0.54
1:O:76:LEU:C	1:O:94:THR:CG2	2.80	0.53
1:B:51:ASN:ND2	1:B:51:ASN:O	2.42	0.53
2:G:23:DA:H2'	2:G:24:DT:C6	2.44	0.53
1:B:78:TYR:HE1	1:B:80:TYR:CD2	2.25	0.53
2:H:15:DC:H2'	2:H:16:DT:C6	2.44	0.53
1:C:50:GLN:HG2	1:C:91:ILE:HD11	1.90	0.53
1:I:41:TYR:C	1:I:41:TYR:CD2	2.86	0.53
1:I:49:THR:HG23	1:I:52:ASP:OD2	2.09	0.53
1:D:127:MET:C	1:D:131:LEU:HD13	2.32	0.53
2:F:20:DT:H2''	2:F:21:DA:C8	2.44	0.53
1:B:82:ASP:O	1:B:85:ASP:O	2.27	0.53
1:C:49:THR:HG21	1:C:88:ARG:HD3	1.91	0.53
1:O:23:LYS:HG3	1:O:107:ILE:HG13	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:22:DA:H2''	2:H:23:DA:H8	1.74	0.52
1:J:39:TYR:CE1	1:J:43:HIS:ND1	2.77	0.52
1:J:106:SER:O	1:J:110:GLU:HG2	2.08	0.52
1:A:24:LEU:HD23	1:A:107:ILE:HD12	1.91	0.52
2:G:15:DC:H2'	2:G:16:DT:C6	2.44	0.52
1:I:112:GLU:OE2	1:J:5:TYR:HB2	2.09	0.52
1:D:41:TYR:CG	1:D:101:VAL:HG11	2.45	0.51
1:B:80:TYR:CE1	1:B:90:ASN:HB2	2.45	0.51
1:C:51:ASN:HB2	2:E:6:DG:OP2	2.10	0.51
2:F:23:DA:H2'	2:F:24:DT:C6	2.45	0.51
1:J:4:THR:O	1:J:8:LEU:HD13	2.10	0.51
1:B:50:GLN:HE22	2:H:6:DG:C3'	2.19	0.51
2:K:1:DA:H2'	2:K:2:DT:H6	1.74	0.51
1:D:54:ALA:HA	1:D:64:VAL:HG21	1.92	0.51
1:O:40:LEU:HD23	1:O:48:LEU:HD12	1.93	0.51
1:C:127:MET:CE	1:D:131:LEU:CD1	2.88	0.50
1:J:101:VAL:HA	1:J:104:PHE:CE2	2.46	0.50
1:J:124:ASN:OD1	1:J:124:ASN:C	2.55	0.50
1:M:49:THR:HB	1:M:52:ASP:H	1.76	0.50
2:L:14:DT:H2'	2:L:15:DC:C6	2.46	0.50
1:J:20:ALA:N	1:J:111:MET:HE1	2.27	0.50
1:M:49:THR:CG2	1:M:51:ASN:H	2.19	0.50
2:H:19:DC:H2'	2:H:20:DT:C6	2.46	0.50
1:C:46:ASP:N	1:C:47:GLY:HA2	2.26	0.50
1:C:33:GLN:HB3	1:C:71:LEU:HD21	1.94	0.50
1:D:31:ASN:OD1	1:D:32:GLU:N	2.45	0.50
2:E:24:DT:H3	2:F:1:DA:H61	1.60	0.50
1:B:79:ARG:HB3	1:B:89:LYS:HB3	1.94	0.49
1:I:127:MET:O	1:I:131:LEU:N	2.35	0.49
2:K:19:DC:H2''	2:K:20:DT:H5'	1.94	0.49
1:C:127:MET:CE	1:D:131:LEU:HD12	2.42	0.49
1:I:59:ARG:NH1	1:J:14:HIS:CD2	2.80	0.49
2:L:15:DC:H2'	2:L:16:DT:C6	2.48	0.49
1:C:106:SER:O	1:C:109:ASP:HB2	2.13	0.49
1:A:89:LYS:NZ	2:G:7:DT:OP1	2.46	0.49
2:H:14:DT:H2'	2:H:15:DC:C6	2.47	0.49
1:C:72:GLU:OE1	1:C:79:ARG:NH2	2.46	0.49
1:M:33:GLN:HE22	1:M:70:ASN:HD22	1.59	0.49
1:I:67:LEU:HD23	1:I:67:LEU:C	2.37	0.49
2:G:10:DG:H2''	2:G:11:DA:OP2	2.13	0.49
2:K:14:DT:H2'	2:K:15:DC:C6	2.48	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:30:THR:OG1	2:L:13:DA:OP1	2.25	0.48
2:H:23:DA:H2''	2:H:24:DT:H5'	1.94	0.48
1:C:101:VAL:O	1:C:105:THR:HB	2.13	0.48
2:G:19:DC:H2'	2:G:20:DT:C6	2.47	0.48
1:M:70:ASN:O	1:M:74:LYS:HG2	2.13	0.48
1:A:101:VAL:HA	1:A:104:PHE:CE2	2.49	0.48
1:C:60:THR:HG23	2:F:15:DC:P	2.54	0.48
1:I:14:HIS:CE1	1:J:59:ARG:HE	2.32	0.48
1:I:46:ASP:N	1:I:47:GLY:HA2	2.28	0.48
2:G:23:DA:H2'	2:G:24:DT:H6	1.79	0.48
1:D:79:ARG:HG2	1:D:91:ILE:CD1	2.43	0.48
2:K:15:DC:H2'	2:K:16:DT:C6	2.49	0.48
1:A:120:SER:O	1:A:124:ASN:ND2	2.44	0.48
1:J:18:GLN:O	1:J:22:GLN:HG3	2.14	0.48
1:A:23:LYS:HD3	1:A:107:ILE:HG23	1.94	0.48
1:C:58:GLN:O	1:C:59:ARG:HD3	2.14	0.48
1:C:112:GLU:O	1:C:116:VAL:HG13	2.13	0.47
1:O:24:LEU:O	1:O:28:ASP:N	2.47	0.47
1:O:103:ALA:O	1:O:107:ILE:HG12	2.14	0.47
1:B:51:ASN:ND2	1:B:51:ASN:C	2.71	0.47
1:D:60:THR:HG23	2:E:15:DC:OP2	2.15	0.47
1:I:16:MET:HE1	1:J:6:SER:HA	1.96	0.47
1:M:41:TYR:HE1	1:M:98:ILE:HG23	1.80	0.47
1:B:118:GLN:C	1:B:119:LEU:HD12	2.39	0.47
1:D:44:GLN:C	1:D:46:ASP:H	2.23	0.47
1:J:43:HIS:O	1:J:46:ASP:HB2	2.14	0.47
1:M:33:GLN:NE2	1:M:71:LEU:HG	2.30	0.47
1:A:112:GLU:O	1:A:116:VAL:HG23	2.14	0.47
2:G:14:DT:H2'	2:G:15:DC:C6	2.50	0.47
1:B:81:VAL:O	1:B:81:VAL:CG1	2.62	0.47
1:C:37:LEU:HD21	1:C:71:LEU:HD13	1.97	0.47
1:M:18:GLN:O	1:M:22:GLN:HG2	2.15	0.47
2:G:22:DA:O3'	2:G:23:DA:H8	1.98	0.47
1:C:127:MET:HE3	1:D:131:LEU:HD12	1.94	0.46
1:J:105:THR:O	1:J:105:THR:HG22	2.16	0.46
1:B:27:PHE:HB3	1:B:100:LEU:HD11	1.96	0.46
1:I:4:THR:HG23	1:I:7:TYR:CB	2.45	0.46
2:E:1:DA:N6	2:F:24:DT:H3	2.13	0.46
1:J:95:THR:O	1:J:99:LYS:HG3	2.16	0.46
1:C:138:LEU:HD21	1:D:115:LEU:HD21	1.96	0.46
1:A:53:ILE:CG2	1:A:57:LEU:HD12	2.45	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:10:ARG:HG2	1:B:11:MET:HE2	1.98	0.46
1:D:101:VAL:HA	1:D:104:PHE:CE2	2.51	0.46
1:A:135:LEU:HD21	1:B:11:MET:HB3	1.98	0.45
1:C:16:MET:HE2	1:C:16:MET:HB3	1.86	0.45
2:F:14:DT:H2'	2:F:15:DC:C6	2.52	0.45
2:L:6:DG:C8	2:L:7:DT:H72	2.52	0.45
2:G:2:DT:H2''	2:G:3:DT:H71	1.98	0.45
1:B:47:GLY:O	1:B:90:ASN:ND2	2.50	0.45
1:O:106:SER:O	1:O:110:GLU:N	2.43	0.45
1:C:87:ARG:HD3	2:F:22:DA:N3	2.32	0.45
1:O:60:THR:O	1:O:64:VAL:HG23	2.17	0.45
1:A:54:ALA:HB1	1:A:59:ARG:O	2.17	0.45
1:B:4:THR:HG22	1:B:6:SER:N	2.27	0.45
2:K:23:DA:H2'	2:K:24:DT:C6	2.51	0.45
1:A:5:TYR:HD2	1:B:128:LYS:HZ3	1.55	0.45
1:O:49:THR:H	1:O:52:ASP:HB2	1.82	0.45
1:J:21:ASP:O	1:J:25:GLU:N	2.49	0.45
1:D:97:GLY:O	1:D:101:VAL:HG23	2.17	0.44
1:J:51:ASN:ND2	2:L:6:DG:OP2	2.50	0.44
1:O:94:THR:HG23	1:O:97:GLY:H	1.82	0.44
2:F:15:DC:H2'	2:F:16:DT:C6	2.52	0.44
1:A:60:THR:HG23	2:H:15:DC:P	2.56	0.44
1:I:127:MET:O	1:I:131:LEU:CB	2.65	0.44
2:H:2:DT:H2''	2:H:3:DT:C6	2.51	0.44
1:D:130:ASN:O	1:D:134:MET:HG3	2.18	0.44
1:I:4:THR:HG23	1:I:7:TYR:HB2	1.99	0.44
1:M:82:ASP:HB2	1:M:90:ASN:ND2	2.32	0.44
2:K:20:DT:OP2	2:K:20:DT:H2'	2.18	0.44
1:A:85:ASP:HB3	1:A:88:ARG:HG3	2.00	0.44
1:M:125:GLU:HA	1:M:128:LYS:HB3	1.99	0.44
1:B:60:THR:HG23	2:G:15:DC:P	2.57	0.44
1:C:30:THR:OG1	1:C:33:GLN:HG3	2.17	0.44
1:I:101:VAL:HA	1:I:104:PHE:CE2	2.53	0.44
1:J:122:GLU:O	1:J:125:GLU:CB	2.65	0.44
1:M:49:THR:CG2	1:M:50:GLN:H	2.27	0.44
1:A:46:ASP:N	1:A:47:GLY:HA2	2.33	0.43
1:B:78:TYR:CE1	1:B:80:TYR:HD2	2.34	0.43
1:M:41:TYR:HA	1:M:93:LEU:HD22	2.00	0.43
1:O:80:TYR:O	1:O:90:ASN:N	2.47	0.43
1:J:10:ARG:HB3	1:J:11:MET:HE2	2.00	0.43
1:M:93:LEU:HD23	1:M:98:ILE:HG12	1.99	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:60:THR:HG23	2:H:15:DC:OP2	2.18	0.43
1:B:115:LEU:HD23	1:B:115:LEU:HA	1.86	0.43
1:D:79:ARG:HG2	1:D:91:ILE:HD12	2.00	0.43
1:M:41:TYR:CE1	1:M:98:ILE:HG23	2.53	0.43
1:D:49:THR:HG21	1:D:88:ARG:HD3	2.01	0.43
1:D:100:LEU:HG	1:D:104:PHE:CD2	2.54	0.43
2:L:22:DA:H2''	2:L:23:DA:H8	1.84	0.43
1:O:106:SER:OG	1:O:107:ILE:N	2.51	0.43
1:B:79:ARG:HA	1:B:90:ASN:O	2.19	0.42
1:M:130:ASN:O	1:M:134:MET:HG2	2.19	0.42
1:A:21:ASP:OD1	1:A:31:ASN:CB	2.67	0.42
1:A:33:GLN:O	1:A:37:LEU:HD12	2.18	0.42
1:A:58:GLN:NE2	1:A:58:GLN:HA	2.35	0.42
1:M:22:GLN:O	1:M:26:GLN:HG3	2.20	0.42
1:C:75:LYS:HE3	1:C:75:LYS:HB2	1.75	0.42
2:L:23:DA:H2'	2:L:24:DT:C6	2.54	0.42
1:O:54:ALA:HB2	1:O:64:VAL:HG21	2.01	0.42
2:F:23:DA:H2'	2:F:24:DT:H6	1.82	0.42
1:A:122:GLU:O	1:A:126:GLN:HG2	2.20	0.42
1:B:101:VAL:HA	1:B:104:PHE:CE2	2.55	0.42
1:C:85:ASP:OD1	2:F:23:DA:H5'	2.19	0.42
1:D:49:THR:CG2	1:D:88:ARG:HD3	2.50	0.42
1:D:4:THR:HG22	1:D:6:SER:H	1.85	0.42
1:M:80:TYR:CZ	1:M:90:ASN:HB2	2.55	0.41
1:A:27:PHE:O	1:A:29:ILE:HG23	2.21	0.41
1:D:112:GLU:O	1:D:116:VAL:HG23	2.19	0.41
1:I:60:THR:HG23	2:L:15:DC:P	2.60	0.41
1:I:127:MET:HA	1:J:127:MET:CE	2.50	0.41
1:J:134:MET:O	1:J:137:SER:HB2	2.20	0.41
2:E:3:DT:H1'	2:E:4:DT:H5'	2.02	0.41
2:E:22:DA:H2''	2:E:23:DA:C8	2.54	0.41
2:F:7:DT:H5'	2:F:7:DT:H6	1.84	0.41
1:I:44:GLN:OE1	1:I:93:LEU:N	2.47	0.41
1:M:82:ASP:OD2	1:M:88:ARG:NH2	2.52	0.41
1:D:43:HIS:CB	1:D:48:LEU:HD11	2.50	0.41
1:O:45:GLN:H	1:O:45:GLN:CD	2.28	0.41
1:C:23:LYS:HD3	1:C:107:ILE:HG23	2.02	0.41
1:D:50:GLN:O	1:D:53:ILE:HB	2.21	0.41
1:D:58:GLN:O	1:D:59:ARG:HD2	2.20	0.41
1:I:130:ASN:CB	1:J:127:MET:HE1	2.51	0.41
1:M:30:THR:HB	1:M:33:GLN:HG3	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:87:ARG:NH1	2:G:4:DT:C1'	2.80	0.41
1:J:78:TYR:CZ	1:J:92:GLY:HA3	2.56	0.41
1:O:49:THR:HG22	1:O:50:GLN:N	2.36	0.41
1:I:50:GLN:NE2	2:K:6:DG:H3'	2.35	0.41
1:D:27:PHE:HB3	1:D:100:LEU:CD1	2.43	0.41
1:D:28:ASP:OD2	1:D:74:LYS:HE2	2.21	0.41
1:D:37:LEU:HD21	1:D:71:LEU:HD22	2.02	0.41
1:I:32:GLU:HG2	2:L:13:DA:O3'	2.21	0.41
1:J:37:LEU:O	1:J:40:LEU:HB2	2.20	0.41
1:J:59:ARG:HD3	1:J:59:ARG:HA	1.81	0.41
1:C:37:LEU:CD2	1:C:71:LEU:HD13	2.51	0.40
1:O:71:LEU:CB	1:O:77:ILE:HG12	2.52	0.40
1:B:45:GLN:CD	1:B:45:GLN:H	2.29	0.40
1:B:86:THR:HB	2:G:22:DA:H4'	2.03	0.40
1:C:107:ILE:O	1:C:111:MET:HG3	2.21	0.40
1:D:79:ARG:NH2	2:F:7:DT:OP1	2.54	0.40
1:J:60:THR:O	1:J:64:VAL:HG23	2.21	0.40
1:M:71:LEU:HD23	1:M:71:LEU:HA	1.98	0.40
1:C:54:ALA:HA	1:C:64:VAL:HG21	2.03	0.40
1:B:48:LEU:HD23	1:B:48:LEU:HA	1.63	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:O:79:ARG:NH2	2:E:22:DA:OP1[1_655]	1.48	0.72
1:C:45:GLN:O	1:D:96:SER:N[2_765]	2.08	0.12

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	136/140 (97%)	132 (97%)	4 (3%)	0	100	100
1	B	136/140 (97%)	131 (96%)	5 (4%)	0	100	100
1	C	136/140 (97%)	135 (99%)	1 (1%)	0	100	100
1	D	135/140 (96%)	133 (98%)	2 (2%)	0	100	100
1	I	135/140 (96%)	129 (96%)	6 (4%)	0	100	100
1	J	136/140 (97%)	131 (96%)	5 (4%)	0	100	100
1	M	120/140 (86%)	119 (99%)	1 (1%)	0	100	100
1	O	133/140 (95%)	130 (98%)	3 (2%)	0	100	100
All	All	1067/1120 (95%)	1040 (98%)	27 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	112/127 (88%)	111 (99%)	1 (1%)	70	82
1	B	115/127 (91%)	113 (98%)	2 (2%)	53	75
1	C	105/127 (83%)	103 (98%)	2 (2%)	50	73
1	D	102/127 (80%)	98 (96%)	4 (4%)	28	59
1	I	87/127 (68%)	84 (97%)	3 (3%)	32	63
1	J	94/127 (74%)	93 (99%)	1 (1%)	65	80
1	M	80/127 (63%)	77 (96%)	3 (4%)	29	60
1	O	80/127 (63%)	76 (95%)	4 (5%)	22	53
All	All	775/1016 (76%)	755 (97%)	20 (3%)	40	69

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	73	ARG
1	B	46	ASP

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Mol	Chain	Res	Type
1	B	99	LYS
1	C	53	ILE
1	C	67	LEU
1	D	3	PHE
1	D	66	ASN
1	D	67	LEU
1	D	127	MET
1	I	18	GLN
1	I	49	THR
1	I	81	VAL
1	J	58	GLN
1	M	31	ASN
1	M	84	GLN
1	M	95	THR
1	O	79	ARG
1	O	81	VAL
1	O	86	THR
1	O	96	SER

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

Mol	Chain	Res	Type
1	A	43	HIS
1	A	50	GLN
1	A	130	ASN
1	B	50	GLN
1	B	51	ASN
1	B	66	ASN
1	C	18	GLN
1	C	58	GLN
1	D	50	GLN
1	D	51	ASN
1	I	58	GLN
1	J	18	GLN
1	J	31	ASN
1	J	33	GLN
1	J	113	GLN
1	M	33	GLN
1	M	44	GLN
1	M	66	ASN
1	M	70	ASN
1	M	90	ASN

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Mol	Chain	Res	Type
1	O	18	GLN
1	O	35	HIS
1	O	66	ASN
1	O	70	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

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	138/140 (98%)	-0.54	0 100 100	26, 45, 101, 135	2 (1%)
1	B	138/140 (98%)	-0.50	1 (0%) 84 66	36, 57, 124, 174	2 (1%)
1	C	138/140 (98%)	-0.52	0 100 100	32, 51, 132, 221	2 (1%)
1	D	137/140 (97%)	-0.22	2 (1%) 72 48	39, 66, 113, 184	2 (1%)
1	I	137/140 (97%)	-0.30	2 (1%) 72 48	50, 77, 126, 163	2 (1%)
1	J	138/140 (98%)	-0.29	3 (2%) 62 38	55, 90, 152, 169	2 (1%)
1	M	124/140 (88%)	-0.35	1 (0%) 82 63	44, 77, 154, 202	1 (0%)
1	O	135/140 (96%)	-0.23	2 (1%) 72 48	47, 93, 184, 226	1 (0%)
2	E	24/24 (100%)	-0.52	0 100 100	40, 56, 102, 116	0
2	F	24/24 (100%)	-0.78	0 100 100	38, 62, 88, 108	0
2	G	24/24 (100%)	-0.64	0 100 100	31, 51, 96, 112	0
2	H	24/24 (100%)	-0.70	0 100 100	27, 57, 101, 119	0
2	K	24/24 (100%)	-0.72	0 100 100	69, 82, 123, 127	0
2	L	24/24 (100%)	-0.61	0 100 100	63, 82, 105, 125	0
All	All	1229/1264 (97%)	-0.40	11 (0%) 81 60	26, 72, 145, 226	14 (1%)

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	O	51	ASN	5.5
1	D	47	GLY	4.4
1	J	87	ARG	3.0
1	O	65	SER	3.0
1	I	16	MET	3.0
1	D	93	LEU	2.8
1	J	9	PHE	2.7
1	B	111	MET	2.4

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Mol	Chain	Res	Type	RSRZ
1	I	80	TYR	2.3
1	J	66	ASN	2.2
1	M	125	GLU	2.1

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no 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.