



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 01:17 pm GMT

PDB ID : 4S04
Title : Crystal structure of *Klebsiella pneumoniae* PmrA in complex with PmrA box DNA
Authors : Hsiao, C.D.; Weng, T.H.; Li, Y.C.
Deposited on : 2014-12-30
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

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : trunk28620
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : recalc28949

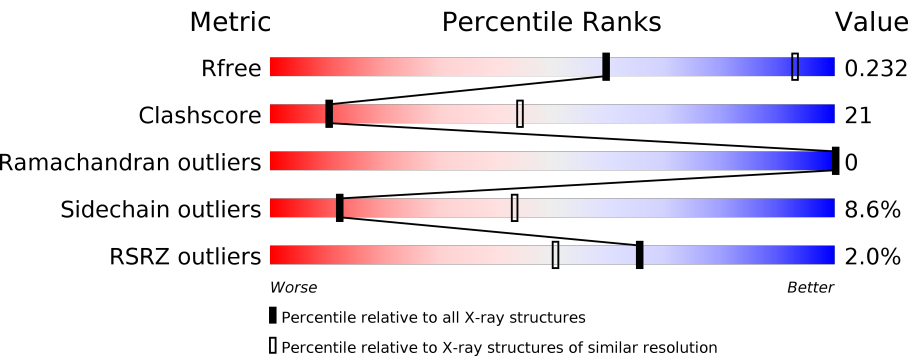
1 Overall quality at a glance i

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}	100719	1015 (3.22-3.18)
Clashscore	112137	1009 (3.20-3.20)
Ramachandran outliers	110173	1118 (3.22-3.18)
Sidechain outliers	110143	1117 (3.22-3.18)
RSRZ outliers	101464	1020 (3.22-3.18)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	232	<div><div></div><div>60%30%6%</div></div>
1	B	232	<div><div>5%</div><div>61%31%6%</div></div>
1	E	232	<div><div>%</div><div>58%29%7%6%</div></div>
1	F	232	<div><div>3%</div><div>58%34%6%</div></div>
2	C	25	<div><div>16%48%36%</div></div>
2	G	25	<div><div>8%24%68%</div></div>

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Mol	Chain	Length	Quality of chain
3	D	25	
3	H	25	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	BEF	E	301	-	-	-	X

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 9060 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 DNA-binding transcriptional regulator BasR.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	219	Total	C	N	O	S	0	0	0
			1726	1080	312	326	8			
1	B	219	Total	C	N	O	S	0	0	0
			1726	1080	312	326	8			
1	E	219	Total	C	N	O	S	0	0	0
			1726	1080	312	326	8			
1	F	219	Total	C	N	O	S	0	0	0
			1726	1080	312	326	8			

There are 44 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	181	GLY	TRP	ENGINEERED MUTATION	UNP S5YJU7
A	220	ASP	ILE	ENGINEERED MUTATION	UNP S5YJU7
A	224	HIS	-	EXPRESSION TAG	UNP S5YJU7
A	225	LEU	-	EXPRESSION TAG	UNP S5YJU7
A	226	GLU	-	EXPRESSION TAG	UNP S5YJU7
A	227	HIS	-	EXPRESSION TAG	UNP S5YJU7
A	228	HIS	-	EXPRESSION TAG	UNP S5YJU7
A	229	HIS	-	EXPRESSION TAG	UNP S5YJU7
A	230	HIS	-	EXPRESSION TAG	UNP S5YJU7
A	231	HIS	-	EXPRESSION TAG	UNP S5YJU7
A	232	HIS	-	EXPRESSION TAG	UNP S5YJU7
B	181	GLY	TRP	ENGINEERED MUTATION	UNP S5YJU7
B	220	ASP	ILE	ENGINEERED MUTATION	UNP S5YJU7
B	224	HIS	-	EXPRESSION TAG	UNP S5YJU7
B	225	LEU	-	EXPRESSION TAG	UNP S5YJU7
B	226	GLU	-	EXPRESSION TAG	UNP S5YJU7
B	227	HIS	-	EXPRESSION TAG	UNP S5YJU7
B	228	HIS	-	EXPRESSION TAG	UNP S5YJU7
B	229	HIS	-	EXPRESSION TAG	UNP S5YJU7
B	230	HIS	-	EXPRESSION TAG	UNP S5YJU7
B	231	HIS	-	EXPRESSION TAG	UNP S5YJU7

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Chain	Residue	Modelled	Actual	Comment	Reference
B	232	HIS	-	EXPRESSION TAG	UNP S5YJU7
E	181	GLY	TRP	ENGINEERED MUTATION	UNP S5YJU7
E	220	ASP	ILE	ENGINEERED MUTATION	UNP S5YJU7
E	224	HIS	-	EXPRESSION TAG	UNP S5YJU7
E	225	LEU	-	EXPRESSION TAG	UNP S5YJU7
E	226	GLU	-	EXPRESSION TAG	UNP S5YJU7
E	227	HIS	-	EXPRESSION TAG	UNP S5YJU7
E	228	HIS	-	EXPRESSION TAG	UNP S5YJU7
E	229	HIS	-	EXPRESSION TAG	UNP S5YJU7
E	230	HIS	-	EXPRESSION TAG	UNP S5YJU7
E	231	HIS	-	EXPRESSION TAG	UNP S5YJU7
E	232	HIS	-	EXPRESSION TAG	UNP S5YJU7
F	181	GLY	TRP	ENGINEERED MUTATION	UNP S5YJU7
F	220	ASP	ILE	ENGINEERED MUTATION	UNP S5YJU7
F	224	HIS	-	EXPRESSION TAG	UNP S5YJU7
F	225	LEU	-	EXPRESSION TAG	UNP S5YJU7
F	226	GLU	-	EXPRESSION TAG	UNP S5YJU7
F	227	HIS	-	EXPRESSION TAG	UNP S5YJU7
F	228	HIS	-	EXPRESSION TAG	UNP S5YJU7
F	229	HIS	-	EXPRESSION TAG	UNP S5YJU7
F	230	HIS	-	EXPRESSION TAG	UNP S5YJU7
F	231	HIS	-	EXPRESSION TAG	UNP S5YJU7
F	232	HIS	-	EXPRESSION TAG	UNP S5YJU7

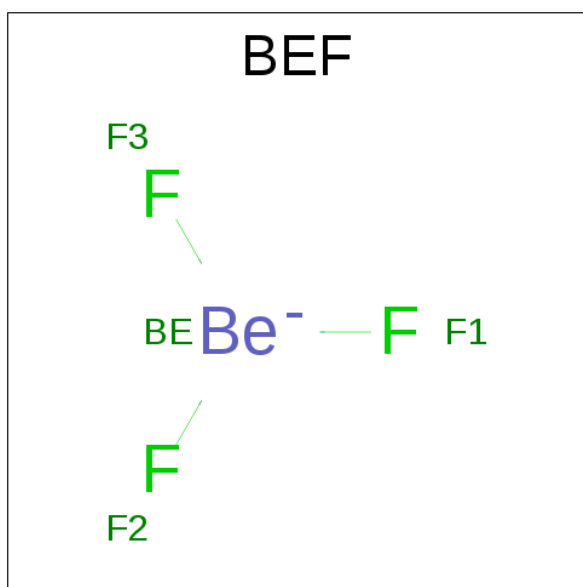
- Molecule 2 is a DNA chain called DNA (25-MER).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	25	Total	C	N	O	P	1	0	0
			509	246	87	151	25			
2	G	25	Total	C	N	O	P	2	0	0
			509	246	87	151	25			

- Molecule 3 is a DNA chain called DNA (25-MER).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	25	Total	C	N	O	P	5	0	0
			516	248	94	149	25			
3	H	25	Total	C	N	O	P	0	0	0
			516	248	94	149	25			

- Molecule 4 is BERYLLIUM TRIFLUORIDE ION (three-letter code: BEF) (formula: BeF₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	Be	F	0	0
			4	1	3		
4	B	1	Total	Be	F	0	0
			4	1	3		
4	E	1	Total	Be	F	0	0
			4	1	3		
4	F	1	Total	Be	F	0	0
			4	1	3		

- Molecule 5 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	1	Total	Mg	0	0
			1	1		
5	A	1	Total	Mg	0	0
			1	1		
5	F	1	Total	Mg	0	0
			1	1		
5	E	1	Total	Mg	0	0
			1	1		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	15	Total	O	0	0
			15	15		

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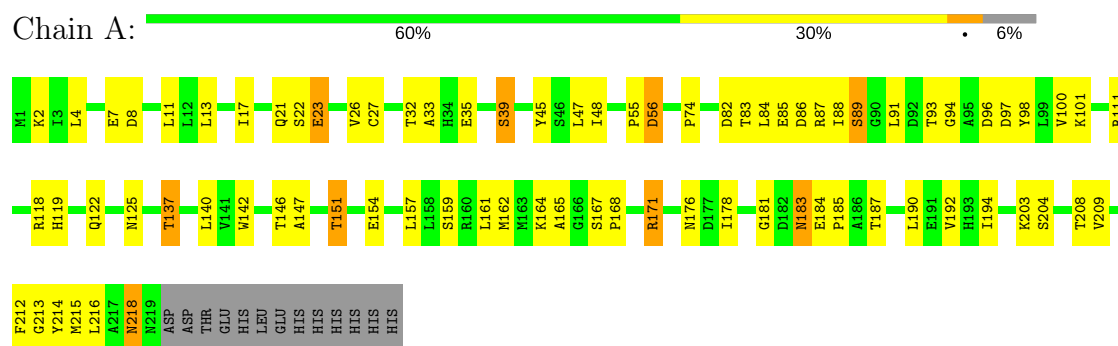
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	13	Total 13	O 13	0	0
6	C	17	Total 17	O 17	0	0
6	D	11	Total 11	O 11	0	0
6	E	11	Total 11	O 11	0	0
6	F	11	Total 11	O 11	0	0
6	G	4	Total 4	O 4	0	0
6	H	4	Total 4	O 4	0	0

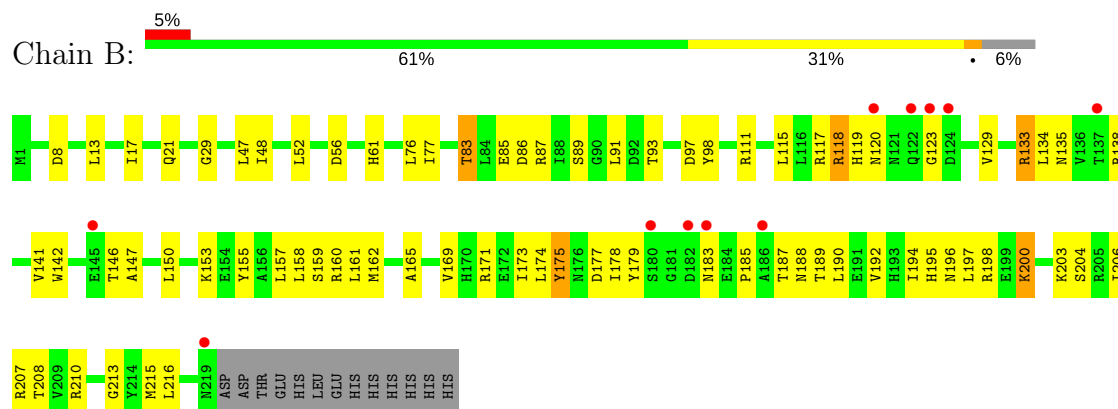
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

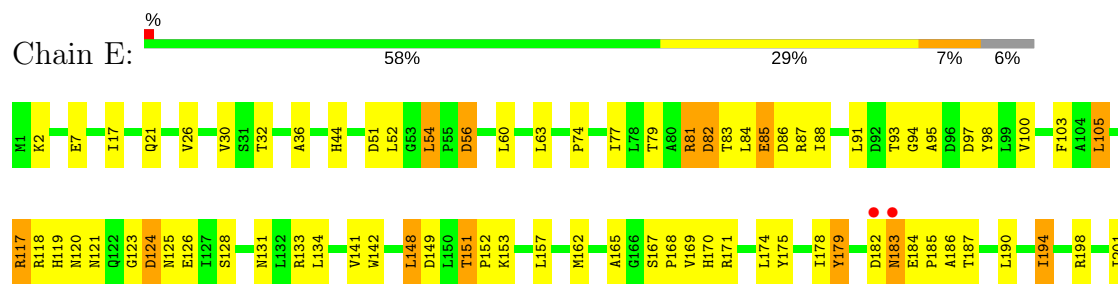
• Molecule 1: DNA-binding transcriptional regulator BasR

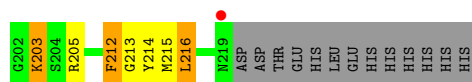


• Molecule 1: DNA-binding transcriptional regulator BasR

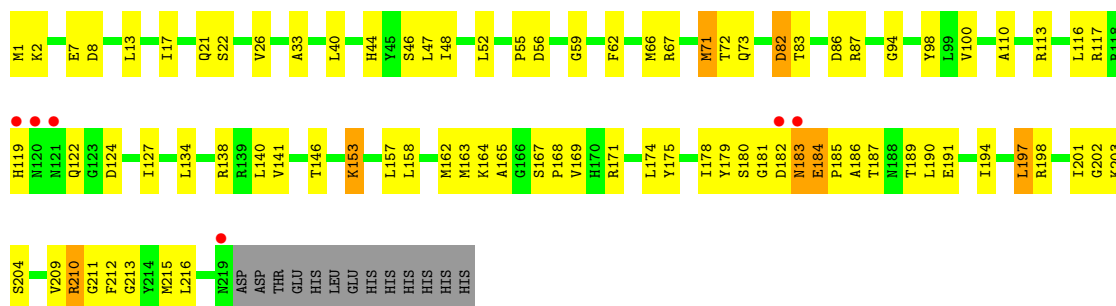


• Molecule 1: DNA-binding transcriptional regulator BasR

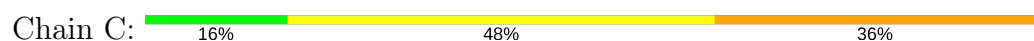




- Molecule 1: DNA-binding transcriptional regulator BasR



- Molecule 2: DNA (25-MER)



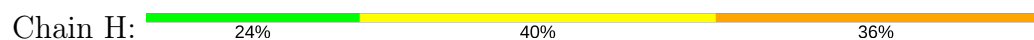
- Molecule 2: DNA (25-MER)



- Molecule 3: DNA (25-MER)



- Molecule 3: DNA (25-MER)



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	194.33Å 250.76Å 108.94Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	24.19 – 3.20 24.19 – 3.20	Depositor EDS
% Data completeness (in resolution range)	45.9 (24.19-3.20) 80.8 (24.19-3.20)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.98 (at 3.23Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.1_1168)	Depositor
R, R_{free}	0.181 , 0.232 0.179 , 0.232	Depositor DCC
R_{free} test set	2022 reflections (5.21%)	DCC
Wilson B-factor (Å ²)	65.2	Xtriage
Anisotropy	0.043	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.26 , 37.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	9060	wwPDB-VP
Average B, all atoms (Å ²)	58.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.22% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

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

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	0.56	1/1750 (0.1%)	0.78	0/2366
1	B	0.50	0/1750	0.78	0/2366
1	E	0.56	0/1750	0.79	0/2366
1	F	0.48	0/1750	0.81	0/2366
2	C	1.08	3/569 (0.5%)	1.83	14/875 (1.6%)
2	G	1.19	6/569 (1.1%)	1.94	22/875 (2.5%)
3	D	1.00	2/579 (0.3%)	1.91	18/892 (2.0%)
3	H	0.97	0/579	2.05	23/892 (2.6%)
All	All	0.70	12/9296 (0.1%)	1.21	77/12998 (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	B	0	1
1	E	0	1
1	F	0	1
All	All	0	3

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	G	16	DC	O3'-P	-8.41	1.51	1.61
2	C	2	DT	O3'-P	-8.01	1.51	1.61
2	G	14	DA	O3'-P	-7.70	1.51	1.61
2	G	3	DT	O3'-P	-7.00	1.52	1.61
2	G	4	DT	O3'-P	-6.38	1.53	1.61
3	D	22	DA	C3'-O3'	-6.04	1.36	1.44
2	G	15	DT	O3'-P	-5.60	1.54	1.61

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	14	DA	C3'-O3'	-5.55	1.36	1.44
3	D	12	DA	C3'-O3'	-5.46	1.36	1.44
1	A	27	CYS	CB-SG	-5.35	1.73	1.81
2	G	23	DA	C3'-O3'	-5.17	1.37	1.44
2	C	1	DA	O3'-P	-5.02	1.55	1.61

All (77) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	H	1	DT	O4'-C1'-N1	-16.43	96.50	108.00
2	G	1	DA	O4'-C1'-N9	-13.52	98.54	108.00
2	C	22	DC	O4'-C1'-N1	-11.23	100.14	108.00
2	G	20	DA	O4'-C1'-N9	11.20	115.84	108.00
2	C	23	DA	O4'-C1'-N9	-10.89	100.38	108.00
2	C	21	DG	O4'-C1'-N9	10.32	115.23	108.00
3	H	4	DC	O4'-C1'-N1	-10.25	100.83	108.00
3	H	24	DT	O4'-C1'-N1	10.19	115.13	108.00
2	G	17	DC	O4'-C1'-N1	-9.22	101.54	108.00
2	C	14	DA	O4'-C1'-N9	-9.09	101.64	108.00
2	G	22	DC	O4'-C1'-N1	9.06	114.34	108.00
2	G	7	DT	O4'-C1'-N1	-8.76	101.86	108.00
3	D	19	DA	O4'-C1'-N9	8.71	114.10	108.00
2	G	10	DT	O4'-C1'-N1	8.62	114.03	108.00
2	G	5	DC	O5'-P-OP2	8.60	121.02	110.70
2	G	25	DG	O4'-C1'-N9	8.50	113.95	108.00
2	G	9	DA	O4'-C1'-N9	8.50	113.95	108.00
3	H	3	DG	O4'-C1'-N9	8.34	113.83	108.00
3	H	6	DT	O4'-C1'-N1	-8.15	102.29	108.00
3	D	23	DA	O4'-C1'-N9	8.14	113.70	108.00
2	C	6	DT	N3-C4-O4	8.03	124.72	119.90
2	G	13	DT	O4'-C4'-C3'	-7.99	101.20	106.00
3	H	13	DA	O4'-C1'-N9	-7.83	102.52	108.00
3	D	5	DT	O4'-C1'-N1	7.82	113.47	108.00
3	D	17	DT	O4'-C1'-N1	-7.74	102.58	108.00
3	D	24	DT	N3-C4-O4	7.63	124.48	119.90
3	H	6	DT	O4'-C4'-C3'	7.55	110.53	106.00
3	H	14	DT	C4-C5-C7	7.50	123.50	119.00
3	D	17	DT	P-O5'-C5'	-7.36	109.13	120.90
2	C	19	DA	O4'-C1'-N9	7.23	113.06	108.00
3	H	6	DT	P-O5'-C5'	-7.04	109.63	120.90
3	D	3	DG	O4'-C4'-C3'	-7.04	101.69	104.50
3	H	3	DG	C1'-O4'-C4'	-6.96	103.14	110.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	H	17	DT	P-O5'-C5'	-6.93	109.81	120.90
2	G	19	DA	O4'-C1'-N9	6.74	112.72	108.00
3	H	14	DT	C6-C5-C7	-6.68	118.89	122.90
3	H	24	DT	P-O5'-C5'	-6.54	110.44	120.90
2	C	15	DT	N3-C4-O4	6.47	123.78	119.90
3	H	17	DT	O4'-C4'-C3'	6.45	109.87	106.00
2	C	13	DT	C1'-O4'-C4'	-6.42	103.68	110.10
2	C	13	DT	O4'-C1'-N1	6.30	112.41	108.00
3	D	24	DT	C5-C4-O4	-6.26	120.52	124.90
3	H	1	DT	N3-C4-O4	6.22	123.63	119.90
2	C	7	DT	P-O5'-C5'	-6.20	110.98	120.90
2	G	16	DC	P-O3'-C3'	-5.92	112.59	119.70
3	H	5	DT	N3-C4-O4	5.88	123.43	119.90
3	D	24	DT	N1-C1'-C2'	-5.88	101.44	112.60
2	G	14	DA	O3'-P-O5'	-5.86	92.87	104.00
2	G	12	DT	N3-C4-O4	5.80	123.38	119.90
2	C	6	DT	C5-C4-O4	-5.79	120.84	124.90
3	D	17	DT	O4'-C4'-C3'	5.78	109.47	106.00
2	G	15	DT	O5'-P-OP2	5.78	117.64	110.70
3	H	1	DT	C5-C4-O4	-5.78	120.86	124.90
3	H	23	DA	O4'-C1'-C2'	5.70	110.46	105.90
3	D	6	DT	N3-C4-O4	5.68	123.31	119.90
3	D	16	DT	N3-C4-O4	5.67	123.31	119.90
2	G	13	DT	C6-C5-C7	-5.67	119.50	122.90
3	H	24	DT	C4'-C3'-C2'	5.60	108.14	103.10
3	D	9	DG	O4'-C1'-N9	5.56	111.89	108.00
2	G	14	DA	O4'-C4'-C3'	5.54	109.33	106.00
2	C	20	DA	O4'-C1'-C2'	-5.51	101.49	105.90
2	C	13	DT	N1-C1'-C2'	5.51	123.06	112.60
3	H	1	DT	N3-C2-O2	5.44	125.56	122.30
3	D	3	DG	C1'-O4'-C4'	-5.43	104.67	110.10
3	H	4	DC	P-O5'-C5'	-5.41	112.24	120.90
2	G	3	DT	O4'-C1'-N1	5.38	111.77	108.00
3	D	7	DA	O4'-C1'-N9	5.37	111.76	108.00
2	G	7	DT	O4'-C4'-C3'	5.36	109.22	106.00
2	C	15	DT	C5-C4-O4	-5.32	121.17	124.90
2	G	25	DG	C3'-C2'-C1'	-5.18	96.28	102.50
3	D	7	DA	C3'-C2'-C1'	-5.17	96.30	102.50
3	D	11	DT	N3-C4-O4	5.16	123.00	119.90
3	H	16	DT	O4'-C1'-N1	-5.16	104.39	108.00
2	G	18	DT	P-O5'-C5'	-5.14	112.68	120.90
3	H	2	DT	C4-C5-C7	5.10	122.06	119.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	G	18	DT	O4'-C1'-N1	5.10	111.57	108.00
3	D	14	DT	P-O5'-C5'	-5.06	112.81	120.90

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	123	GLY	Peptide
1	E	124	ASP	Peptide
1	F	184	GLU	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	1726	0	1755	55	0
1	B	1726	0	1755	65	0
1	E	1726	0	1755	90	0
1	F	1726	0	1755	75	0
2	C	509	0	286	22	0
2	G	509	0	286	39	0
3	D	516	0	285	26	0
3	H	516	0	285	22	0
4	A	4	0	0	0	0
4	B	4	0	0	0	0
4	E	4	0	0	0	0
4	F	4	0	0	0	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
5	E	1	0	0	0	0
5	F	1	0	0	0	0
6	A	15	0	0	3	0
6	B	13	0	0	2	0
6	C	17	0	0	1	0
6	D	11	0	0	1	0
6	E	11	0	0	0	0
6	F	11	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	G	4	0	0	3	0
6	H	4	0	0	1	0
All	All	9060	0	8162	352	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 21.

All (352) 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:E:151:THR:HG21	2:G:4:DT:H3'	1.24	1.15
1:E:151:THR:HB	2:G:5:DC:OP2	1.53	1.08
1:E:151:THR:CB	2:G:5:DC:OP2	2.09	1.00
2:G:6:DT:H2''	2:G:7:DT:H5'	1.43	0.98
1:E:178:ILE:HG13	1:E:179:TYR:CE1	2.02	0.95
1:E:179:TYR:CE2	1:E:185:PRO:HG2	2.05	0.91
3:H:16:DT:H2''	3:H:17:DT:H5''	1.54	0.89
1:E:151:THR:CG2	2:G:4:DT:H3'	2.02	0.88
3:D:16:DT:H2''	3:D:17:DT:H5''	1.57	0.87
1:E:151:THR:HG22	2:G:4:DT:H5''	1.56	0.84
1:E:178:ILE:HG13	1:E:179:TYR:CD1	2.13	0.84
1:E:151:THR:CG2	2:G:5:DC:OP2	2.25	0.84
3:H:5:DT:H2''	3:H:6:DT:H5''	1.60	0.83
3:D:5:DT:H2''	3:D:6:DT:H5'	1.58	0.83
1:E:151:THR:CG2	2:G:4:DT:H5''	2.09	0.82
1:F:153:LYS:CG	1:F:178:ILE:HD12	2.11	0.79
1:F:153:LYS:HG2	1:F:178:ILE:HG23	1.64	0.78
1:E:151:THR:HG21	2:G:4:DT:C3'	2.13	0.75
1:F:153:LYS:NZ	1:F:153:LYS:HB2	2.01	0.75
2:G:15:DT:H2''	2:G:16:DC:O5'	1.86	0.75
1:E:178:ILE:HG13	1:E:179:TYR:HE1	1.52	0.74
1:F:153:LYS:HG2	1:F:178:ILE:HD12	1.68	0.73
1:F:157:LEU:HD22	1:F:174:LEU:HD23	1.70	0.73
1:B:161:LEU:HB3	1:B:216:LEU:HD22	1.69	0.73
1:F:171:ARG:HG3	1:F:213:GLY:HA2	1.70	0.72
1:B:187:THR:HG21	2:C:18:DT:H2'	1.71	0.71
1:E:32:THR:HB	1:E:56:ASP:OD2	1.90	0.71
1:E:151:THR:HG21	2:G:5:DC:OP2	1.89	0.71
1:E:179:TYR:CE2	1:E:185:PRO:CG	2.72	0.71
1:F:181:GLY:N	1:F:182:ASP:HA	2.05	0.70
1:F:157:LEU:HG	1:F:178:ILE:HD11	1.74	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:H:16:DT:C2'	3:H:17:DT:H5''	2.21	0.70
1:A:171:ARG:HG3	1:A:213:GLY:HA2	1.73	0.70
1:A:164:LYS:HE3	1:B:138:ARG:HH22	1.56	0.70
1:E:178:ILE:C	1:E:179:TYR:CD1	2.65	0.70
1:F:134:LEU:HD23	1:F:141:VAL:HG22	1.73	0.69
3:H:5:DT:C2'	3:H:6:DT:H5''	2.22	0.69
2:G:3:DT:H2''	2:G:4:DT:O5'	1.93	0.69
2:G:23:DA:H2''	2:G:24:DA:C8	2.28	0.69
1:A:82:ASP:HB3	1:A:100:VAL:HG11	1.75	0.68
3:D:16:DT:C2'	3:D:17:DT:H5''	2.21	0.68
1:E:121:ASN:ND2	1:E:124:ASP:O	2.27	0.67
1:F:180:SER:N	1:F:181:GLY:HA2	2.10	0.67
1:A:74:PRO:HD3	1:A:119:HIS:HD2	1.60	0.66
1:E:179:TYR:HE2	1:E:185:PRO:HG2	1.60	0.66
1:B:153:LYS:HG2	2:C:16:DC:OP1	1.97	0.65
1:E:170:HIS:CD2	1:E:212:PHE:HE1	2.14	0.65
1:A:118:ARG:NH1	1:A:122:GLN:O	2.29	0.65
1:A:74:PRO:HD3	1:A:119:HIS:CD2	2.31	0.65
1:B:97:ASP:OD2	1:B:111:ARG:HD2	1.97	0.64
1:A:157:LEU:HG	1:A:178:ILE:HD11	1.78	0.64
1:E:212:PHE:HE2	1:F:140:LEU:HD13	1.62	0.64
2:G:6:DT:H2''	2:G:7:DT:C5'	2.22	0.64
3:H:1:DT:H4'	3:H:1:DT:OP2	1.98	0.64
3:H:15:DA:H8	3:H:15:DA:H5'	1.62	0.64
2:G:23:DA:H2''	2:G:24:DA:H8	1.62	0.63
1:E:151:THR:CG2	2:G:4:DT:C3'	2.73	0.63
1:F:181:GLY:H	1:F:182:ASP:HA	1.64	0.62
1:A:96:ASP:OD2	1:B:118:ARG:NH2	2.32	0.62
1:E:178:ILE:CG1	1:E:179:TYR:HE1	2.13	0.62
1:A:56:ASP:OD1	1:A:56:ASP:N	2.32	0.62
1:E:51:ASP:O	1:E:54:LEU:HD22	1.99	0.62
1:A:17:ILE:O	1:A:21:GLN:HG3	1.99	0.61
2:G:13:DT:H2'	2:G:14:DA:C8	2.35	0.61
1:B:216:LEU:H	1:B:216:LEU:HD23	1.65	0.61
1:B:153:LYS:HD2	1:B:178:ILE:HG23	1.81	0.61
1:F:1:MET:HB3	1:F:46:SER:HB2	1.82	0.61
1:E:151:THR:HG22	1:E:152:PRO:HD2	1.83	0.61
1:A:181:GLY:HA2	1:A:184:GLU:CD	2.21	0.61
1:B:17:ILE:O	1:B:21:GLN:HG2	2.01	0.61
1:F:153:LYS:HZ3	1:F:153:LYS:HB2	1.64	0.60
1:F:17:ILE:O	1:F:21:GLN:HG3	2.01	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:168:PRO:HA	1:A:215:MET:HB2	1.83	0.60
1:E:178:ILE:CG1	1:E:179:TYR:CE1	2.82	0.60
1:B:158:LEU:HD22	1:B:197:LEU:HD21	1.84	0.60
1:F:153:LYS:HG3	1:F:178:ILE:HD12	1.80	0.60
2:G:4:DT:H1'	6:G:102:HOH:O	2.00	0.60
1:F:153:LYS:CG	1:F:178:ILE:HG23	2.32	0.60
2:G:19:DA:H1'	6:G:103:HOH:O	2.01	0.59
1:A:218:ASN:N	1:A:218:ASN:OD1	2.34	0.59
1:E:17:ILE:O	1:E:21:GLN:HG3	2.01	0.59
1:F:203:LYS:N	1:F:204:SER:HA	2.18	0.59
1:A:23:GLU:HG2	6:A:401:HOH:O	2.03	0.59
1:F:158:LEU:HB2	1:F:197:LEU:HD21	1.84	0.59
1:B:133:ARG:HG3	1:B:142:TRP:HB2	1.85	0.58
1:B:203:LYS:HG3	1:B:204:SER:HB3	1.84	0.58
1:E:124:ASP:HB3	1:F:67:ARG:HD2	1.84	0.58
1:F:198:ARG:O	1:F:202:GLY:HA3	2.03	0.58
1:E:51:ASP:O	1:E:54:LEU:CD2	2.52	0.58
1:F:171:ARG:NH2	1:F:191:GLU:OE2	2.35	0.58
1:E:178:ILE:C	1:E:179:TYR:HD1	2.04	0.58
1:F:210:ARG:NE	3:H:2:DT:H4'	2.19	0.58
1:A:35:GLU:O	1:A:39:SER:OG	2.20	0.58
1:E:83:THR:HB	1:E:86:ASP:H	1.68	0.57
1:A:82:ASP:O	1:A:87:ARG:NH2	2.37	0.57
1:E:190:LEU:HD11	1:E:194:ILE:HD11	1.87	0.57
1:F:82:ASP:OD1	1:F:82:ASP:N	2.37	0.57
1:E:123:GLY:HA2	1:F:72:THR:HB	1.84	0.57
1:E:83:THR:HG22	1:E:85:GLU:H	1.69	0.57
1:B:160:ARG:NH1	1:B:173:ILE:HG22	2.19	0.57
3:H:17:DT:H2''	3:H:18:DA:O4'	2.05	0.57
1:F:169:VAL:HB	1:F:174:LEU:HD11	1.85	0.57
1:F:209:VAL:HG12	1:F:212:PHE:HB3	1.87	0.57
1:F:167:SER:C	1:F:215:MET:HG3	2.25	0.57
1:E:82:ASP:O	1:E:87:ARG:NH2	2.37	0.57
1:B:185:PRO:HB3	2:C:18:DT:OP2	2.05	0.56
1:E:157:LEU:HD21	1:E:178:ILE:HD13	1.86	0.56
2:G:1:DA:H2'	2:G:1:DA:OP2	2.05	0.56
1:B:210:ARG:HG3	3:D:3:DG:H5'	1.86	0.56
1:B:203:LYS:NZ	3:D:5:DT:OP1	2.34	0.56
1:A:171:ARG:HB3	1:A:171:ARG:HH11	1.70	0.56
1:B:160:ARG:HH12	1:B:173:ILE:HG22	1.70	0.56
1:E:7:GLU:OE2	1:E:51:ASP:HB2	2.06	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:186:ALA:HB3	2:G:18:DT:OP2	2.06	0.56
1:A:151:THR:HB	2:C:4:DT:H3'	1.87	0.55
1:E:179:TYR:CD1	1:E:179:TYR:N	2.73	0.55
1:E:79:THR:HG23	1:E:81:ARG:H	1.70	0.55
3:H:8:DG:H2''	3:H:9:DG:C8	2.42	0.55
3:H:15:DA:C8	3:H:15:DA:H5'	2.40	0.55
1:A:97:ASP:OD2	1:A:111:ARG:HD2	2.06	0.54
1:B:13:LEU:HD21	1:B:29:GLY:HA3	1.89	0.54
1:F:187:THR:HG21	2:G:18:DT:H2'	1.89	0.54
1:F:33:ALA:HB3	1:F:56:ASP:OD2	2.07	0.54
1:A:151:THR:HG23	1:A:154:GLU:CD	2.27	0.54
2:C:23:DA:H2''	2:C:24:DA:C8	2.43	0.54
1:E:198:ARG:HD3	1:E:203:LYS:HB2	1.88	0.54
3:H:9:DG:OP2	3:H:9:DG:H2'	2.07	0.54
3:D:15:DA:H5'	3:D:15:DA:H8	1.72	0.54
1:E:84:LEU:O	1:E:88:ILE:HG12	2.08	0.53
1:B:61:HIS:HB2	6:B:412:HOH:O	2.08	0.53
1:E:179:TYR:CD2	1:E:185:PRO:CG	2.92	0.53
1:E:198:ARG:HA	1:E:201:ILE:HG22	1.89	0.53
1:A:171:ARG:NH1	1:A:171:ARG:HB3	2.23	0.53
1:E:169:VAL:O	1:E:213:GLY:HA3	2.09	0.53
2:C:22:DC:H2''	2:C:23:DA:C8	2.43	0.53
1:A:97:ASP:OD1	1:A:98:TYR:N	2.39	0.52
1:E:167:SER:OG	1:F:138:ARG:NH2	2.42	0.52
1:B:204:SER:HB2	1:B:206:ILE:O	2.09	0.52
1:B:158:LEU:HB2	1:B:197:LEU:HD11	1.91	0.52
2:C:13:DT:H2'	2:C:14:DA:N7	2.24	0.52
1:B:157:LEU:HD22	1:B:174:LEU:HD13	1.92	0.52
3:H:1:DT:H2'	3:H:2:DT:C6	2.45	0.52
1:A:33:ALA:HB3	1:A:56:ASP:OD2	2.10	0.52
2:G:21:DG:H2''	2:G:22:DC:OP2	2.10	0.52
1:B:135:ASN:HB3	1:B:138:ARG:H	1.75	0.51
3:H:2:DT:H2''	3:H:3:DG:O4'	2.11	0.51
1:E:153:LYS:HD2	1:E:178:ILE:HB	1.92	0.51
1:F:2:LYS:HG3	1:F:26:VAL:HG12	1.92	0.51
1:E:168:PRO:HB2	1:E:212:PHE:CD2	2.45	0.51
1:F:134:LEU:HD23	1:F:141:VAL:CG2	2.39	0.51
3:D:7:DA:H2''	3:D:8:DG:N7	2.25	0.51
1:B:198:ARG:HH22	1:B:208:THR:HB	1.76	0.51
1:E:151:THR:HB	2:G:5:DC:P	2.50	0.51
1:B:150:LEU:HD11	1:B:158:LEU:HD23	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:185:PRO:HG2	1:B:189:THR:HG21	1.92	0.50
1:A:203:LYS:HA	6:A:413:HOH:O	2.11	0.50
1:A:8:ASP:HB3	1:A:55:PRO:HD3	1.93	0.50
1:B:160:ARG:HG2	1:B:160:ARG:HH11	1.77	0.50
1:B:159:SER:OG	1:B:160:ARG:N	2.45	0.50
2:C:12:DT:H2''	2:C:13:DT:C6	2.45	0.50
2:C:13:DT:H2'	2:C:14:DA:C8	2.47	0.50
1:E:162:MET:HA	1:E:165:ALA:HB2	1.93	0.50
1:A:171:ARG:CB	1:A:171:ARG:HH11	2.23	0.50
1:A:190:LEU:O	1:A:194:ILE:HG12	2.12	0.50
1:E:126:GLU:HB2	1:E:134:LEU:O	2.11	0.50
2:G:9:DA:H2''	2:G:10:DT:OP2	2.12	0.50
1:B:165:ALA:HA	1:B:216:LEU:HD21	1.92	0.50
1:E:198:ARG:HB3	1:E:203:LYS:HB3	1.93	0.49
1:B:158:LEU:HB2	1:B:197:LEU:HD21	1.94	0.49
2:G:9:DA:OP2	2:G:9:DA:H2'	2.13	0.49
1:A:209:VAL:HG12	1:A:212:PHE:HB3	1.94	0.49
1:A:17:ILE:HG22	1:A:21:GLN:OE1	2.12	0.49
1:A:4:LEU:HB2	1:A:45:TYR:CD2	2.47	0.49
2:C:18:DT:H3	3:D:7:DA:H61	1.60	0.49
2:C:5:DC:H2''	2:C:6:DT:O5'	2.11	0.49
1:A:162:MET:HA	1:A:165:ALA:HB2	1.94	0.49
1:B:134:LEU:HD23	1:B:141:VAL:HG22	1.95	0.49
1:E:178:ILE:CD1	1:E:179:TYR:HE1	2.26	0.49
1:F:179:TYR:CE2	1:F:183:ASN:HB3	2.48	0.49
1:B:83:THR:HB	1:B:86:ASP:H	1.77	0.49
3:D:21:DA:C2	3:D:22:DA:C4	3.00	0.49
1:A:125:ASN:OD1	1:A:137:THR:OG1	2.30	0.49
2:C:21:DG:H2''	6:C:106:HOH:O	2.12	0.49
1:E:2:LYS:HG3	1:E:26:VAL:HG12	1.95	0.49
1:A:183:ASN:OD1	1:A:183:ASN:N	2.44	0.49
1:B:216:LEU:N	1:B:216:LEU:HD23	2.28	0.49
1:E:169:VAL:HB	1:E:174:LEU:HD11	1.94	0.49
1:E:151:THR:CG2	2:G:4:DT:C5'	2.85	0.49
1:F:180:SER:H	1:F:181:GLY:HA2	1.78	0.48
1:E:77:ILE:HG13	1:E:95:ALA:HB2	1.95	0.48
1:B:17:ILE:HD11	1:B:29:GLY:HA3	1.94	0.48
1:A:157:LEU:CG	1:A:178:ILE:HD11	2.42	0.48
1:A:85:GLU:O	1:A:89:SER:OG	2.27	0.48
1:E:118:ARG:HH12	1:E:123:GLY:HA2	1.78	0.48
2:G:1:DA:H8	2:G:1:DA:OP2	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:21:DG:H2''	2:C:22:DC:OP2	2.14	0.48
2:G:24:DA:H2'	2:G:25:DG:C8	2.48	0.48
2:C:21:DG:H1'	2:C:22:DC:H5''	1.94	0.48
3:D:17:DT:H2''	3:D:18:DA:O4'	2.14	0.48
2:G:10:DT:H5'	6:G:104:HOH:O	2.14	0.48
1:E:179:TYR:HE2	1:E:185:PRO:CG	2.21	0.48
1:E:93:THR:N	1:E:94:GLY:CA	2.76	0.48
3:D:11:DT:H2''	3:D:12:DA:C8	2.49	0.48
1:F:87:ARG:HA	1:F:98:TYR:CE2	2.48	0.47
1:E:103:PHE:CE2	1:E:105:LEU:HD12	2.49	0.47
1:E:179:TYR:HD1	1:E:179:TYR:N	2.11	0.47
1:E:74:PRO:HD3	1:E:119:HIS:CD2	2.49	0.47
1:A:7:GLU:OE2	1:A:101:LYS:NZ	2.34	0.47
1:F:164:LYS:HE3	1:F:167:SER:HB2	1.97	0.47
1:A:2:LYS:HA	1:A:26:VAL:O	2.15	0.47
1:F:82:ASP:HB3	1:F:100:VAL:CG1	2.44	0.47
1:F:162:MET:HA	1:F:165:ALA:HB2	1.96	0.47
3:D:18:DA:H2''	3:D:19:DA:O4'	2.15	0.47
1:F:82:ASP:HB3	1:F:100:VAL:HG11	1.96	0.47
1:A:167:SER:C	1:A:215:MET:HG3	2.35	0.47
2:C:14:DA:C2	3:D:12:DA:H2	2.33	0.47
1:B:150:LEU:HB2	1:B:155:TYR:HB2	1.95	0.47
1:B:200:LYS:NZ	1:B:200:LYS:HA	2.30	0.47
2:C:24:DA:H61	3:D:1:DT:H3	1.63	0.47
1:B:160:ARG:NH2	1:B:177:ASP:OD2	2.49	0.46
1:B:215:MET:HG2	1:B:216:LEU:N	2.29	0.46
1:B:83:THR:HG22	1:B:85:GLU:H	1.79	0.46
1:E:175:TYR:OH	1:E:184:GLU:HA	2.15	0.46
1:E:52:LEU:HD12	1:E:81:ARG:HG3	1.96	0.46
1:E:91:LEU:O	1:F:117:ARG:NH1	2.48	0.46
1:B:97:ASP:OD1	1:B:111:ARG:NH2	2.49	0.46
3:D:6:DT:H2''	3:D:7:DA:OP2	2.15	0.46
1:F:52:LEU:HA	1:F:59:GLY:HA3	1.97	0.46
2:G:14:DA:H8	2:G:14:DA:OP2	1.97	0.46
1:A:82:ASP:HB3	1:A:100:VAL:CG1	2.43	0.46
1:E:117:ARG:O	1:E:121:ASN:O	2.33	0.46
3:D:15:DA:C8	3:D:15:DA:H5'	2.50	0.46
1:F:153:LYS:NZ	1:F:153:LYS:CB	2.73	0.46
2:G:12:DT:H2''	2:G:13:DT:C6	2.51	0.46
1:A:142:TRP:CZ3	1:A:147:ALA:HB2	2.50	0.46
1:E:212:PHE:CE2	1:F:140:LEU:HD13	2.46	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:179:TYR:HE2	1:F:183:ASN:HB3	1.80	0.46
1:B:160:ARG:HG2	1:B:160:ARG:NH1	2.31	0.45
1:B:47:LEU:HD23	1:B:48:ILE:N	2.31	0.45
1:E:215:MET:HG2	1:E:216:LEU:N	2.30	0.45
1:A:164:LYS:HE3	1:B:138:ARG:NH2	2.26	0.45
1:B:52:LEU:HD21	1:B:77:ILE:HD13	1.98	0.45
2:C:14:DA:C2	3:D:12:DA:C2	3.04	0.45
1:F:190:LEU:O	1:F:194:ILE:HG12	2.16	0.45
3:D:18:DA:H8	6:D:103:HOH:O	1.98	0.45
1:F:153:LYS:HZ2	1:F:153:LYS:HB2	1.82	0.45
1:A:203:LYS:HA	1:A:204:SER:HA	1.49	0.45
1:A:47:LEU:HD23	1:A:48:ILE:N	2.32	0.45
1:B:146:THR:HG22	1:B:147:ALA:H	1.82	0.45
1:B:198:ARG:HB3	1:B:203:LYS:HB2	1.99	0.45
1:F:215:MET:HG2	1:F:216:LEU:N	2.32	0.44
2:C:19:DA:H2''	2:C:20:DA:C8	2.52	0.44
1:B:215:MET:HG2	1:B:216:LEU:H	1.82	0.44
1:B:89:SER:O	1:B:93:THR:OG1	2.25	0.44
1:E:131:ASN:CG	1:E:205:ARG:HH12	2.20	0.44
1:E:133:ARG:HB3	1:E:142:TRP:HB2	1.99	0.44
1:E:141:VAL:CG1	1:E:148:LEU:HD11	2.48	0.44
1:A:184:GLU:HA	1:A:185:PRO:HD3	1.86	0.44
1:A:215:MET:HG2	1:A:216:LEU:N	2.33	0.44
2:C:24:DA:H2'	2:C:25:DG:N7	2.33	0.44
1:F:127:ILE:HD12	1:F:163:MET:CE	2.48	0.44
3:H:11:DT:H2''	3:H:12:DA:C8	2.53	0.44
3:D:11:DT:H2'	3:D:11:DT:H6	1.46	0.44
1:F:8:ASP:N	1:F:8:ASP:OD1	2.44	0.44
1:B:183:ASN:OD1	1:B:183:ASN:N	2.46	0.44
1:B:190:LEU:O	1:B:194:ILE:HG12	2.17	0.44
2:C:3:DT:O2	3:D:23:DA:H2	2.00	0.44
1:E:103:PHE:HE2	1:E:105:LEU:HD12	1.83	0.44
1:F:153:LYS:HB2	2:G:16:DC:OP1	2.17	0.44
1:A:161:LEU:HD23	1:A:161:LEU:HA	1.78	0.43
1:B:56:ASP:N	1:B:56:ASP:OD1	2.44	0.43
1:E:133:ARG:NH1	1:E:142:TRP:CZ3	2.76	0.43
1:E:97:ASP:OD1	1:E:98:TYR:N	2.51	0.43
1:F:157:LEU:HD21	1:F:178:ILE:HG12	2.00	0.43
1:F:87:ARG:HA	1:F:98:TYR:CD2	2.53	0.43
2:G:4:DT:O2	3:H:22:DA:C2	2.71	0.43
1:B:162:MET:HA	1:B:165:ALA:HB2	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:175:TYR:O	1:B:179:TYR:HD2	2.02	0.43
1:E:215:MET:HG2	1:E:216:LEU:H	1.84	0.43
1:F:168:PRO:HB2	1:F:212:PHE:CE2	2.53	0.43
2:G:7:DT:H2''	2:G:8:DA:H5'	2.00	0.43
1:E:63:LEU:HA	1:E:63:LEU:HD12	1.72	0.43
1:F:40:LEU:HD13	1:F:71:MET:HE1	1.99	0.43
3:D:23:DA:H2''	3:D:24:DT:O5'	2.19	0.43
2:G:24:DA:C6	2:G:25:DG:C6	3.07	0.43
2:G:23:DA:C2	2:G:24:DA:C4	3.07	0.43
3:H:5:DT:C3'	3:H:6:DT:H5''	2.48	0.43
1:A:203:LYS:HG2	6:A:413:HOH:O	2.19	0.43
1:A:7:GLU:HB2	1:A:13:LEU:HD13	2.01	0.43
3:D:24:DT:H2''	3:D:25:DC:O5'	2.19	0.43
2:G:18:DT:H2''	2:G:19:DA:O4'	2.19	0.43
1:B:169:VAL:O	1:B:213:GLY:HA3	2.18	0.42
1:A:214:TYR:HE2	3:D:15:DA:OP1	2.01	0.42
1:A:178:ILE:HD13	1:A:178:ILE:HG21	1.69	0.42
1:B:97:ASP:OD1	1:B:98:TYR:N	2.52	0.42
1:F:201:ILE:HB	1:F:202:GLY:CA	2.50	0.42
1:F:33:ALA:HB1	1:F:62:PHE:HD2	1.85	0.42
3:H:15:DA:C8	3:H:16:DT:H72	2.55	0.42
1:E:134:LEU:HD12	1:E:141:VAL:HG22	2.00	0.42
1:E:198:ARG:HD3	1:E:203:LYS:CB	2.49	0.42
1:F:8:ASP:HB3	1:F:55:PRO:HD3	2.01	0.42
1:B:91:LEU:HA	1:B:91:LEU:HD23	1.75	0.42
1:E:84:LEU:HD11	1:E:88:ILE:HD11	2.02	0.42
1:E:88:ILE:HG23	1:F:110:ALA:HB2	2.00	0.42
1:A:84:LEU:O	1:A:88:ILE:HG13	2.19	0.42
1:B:196:ASN:HA	6:B:401:HOH:O	2.18	0.42
1:E:183:ASN:O	1:E:184:GLU:C	2.58	0.42
1:A:83:THR:OG1	1:A:86:ASP:OD2	2.35	0.42
1:B:118:ARG:HD2	1:B:118:ARG:HA	1.87	0.42
1:F:47:LEU:C	1:F:48:ILE:HD12	2.40	0.42
1:B:129:VAL:HG11	1:B:165:ALA:CB	2.50	0.42
1:B:188:ASN:O	1:B:192:VAL:HG23	2.20	0.42
1:E:178:ILE:O	1:E:179:TYR:CD1	2.73	0.42
3:H:22:DA:H2''	3:H:23:DA:O4'	2.19	0.42
1:E:214:TYR:N	1:E:214:TYR:CD1	2.87	0.41
1:F:179:TYR:CE1	1:F:189:THR:HG21	2.55	0.41
3:H:8:DG:H2''	3:H:9:DG:H8	1.84	0.41
1:E:117:ARG:CZ	1:F:94:GLY:HA2	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:210:ARG:HD2	3:H:2:DT:O3'	2.20	0.41
1:A:32:THR:HB	1:A:56:ASP:OD1	2.21	0.41
1:B:76:LEU:HD23	1:B:77:ILE:O	2.21	0.41
3:D:17:DT:H2'	3:D:18:DA:C8	2.55	0.41
1:F:197:LEU:O	1:F:201:ILE:HG12	2.21	0.41
3:H:14:DT:H2''	3:H:15:DA:OP2	2.21	0.41
1:A:91:LEU:O	1:B:117:ARG:NH1	2.53	0.41
1:B:87:ARG:HD3	1:B:87:ARG:HH21	1.74	0.41
1:E:170:HIS:CD2	1:E:212:PHE:CE1	3.02	0.41
1:F:71:MET:HB3	1:F:71:MET:HE2	1.68	0.41
1:B:111:ARG:O	1:B:115:LEU:HG	2.20	0.41
1:B:8:ASP:N	1:B:8:ASP:OD1	2.52	0.41
2:C:11:DA:C2	2:C:12:DT:C2	3.09	0.41
2:C:6:DT:H2''	2:C:7:DT:H5'	2.01	0.41
1:F:40:LEU:HD23	1:F:40:LEU:HA	1.93	0.41
1:E:117:ARG:NH2	1:E:125:ASN:OD1	2.53	0.41
1:F:204:SER:OG	1:F:204:SER:O	2.36	0.41
1:A:93:THR:HB	1:A:94:GLY:HA2	2.03	0.41
3:D:22:DA:H2'	3:D:23:DA:C8	2.56	0.41
1:E:30:VAL:HG11	1:E:36:ALA:N	2.36	0.40
1:F:175:TYR:CE1	1:F:185:PRO:HG3	2.56	0.40
1:F:67:ARG:HH22	1:F:73:GLN:H	1.68	0.40
1:F:83:THR:HG23	1:F:86:ASP:OD2	2.20	0.40
1:F:116:LEU:O	1:F:119:HIS:HB3	2.21	0.40
1:F:7:GLU:HB3	1:F:13:LEU:HB2	2.03	0.40
3:H:14:DT:H71	6:H:102:HOH:O	2.21	0.40
1:F:167:SER:HA	1:F:168:PRO:HD3	1.99	0.40
1:F:210:ARG:HG3	1:F:211:GLY:N	2.37	0.40
2:G:5:DC:H2''	2:G:6:DT:O5'	2.21	0.40
3:D:11:DT:H4'	3:D:12:DA:OP1	2.21	0.40
1:E:119:HIS:ND1	1:E:120:ASN:ND2	2.69	0.40
1:E:178:ILE:HG13	1:E:179:TYR:HD1	1.75	0.40
1:E:185:PRO:O	1:E:186:ALA:HB3	2.22	0.40
1:E:93:THR:N	1:E:94:GLY:HA2	2.36	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	217/232 (94%)	213 (98%)	4 (2%)	0	100	100
1	B	217/232 (94%)	211 (97%)	6 (3%)	0	100	100
1	E	217/232 (94%)	209 (96%)	8 (4%)	0	100	100
1	F	217/232 (94%)	210 (97%)	7 (3%)	0	100	100
All	All	868/928 (94%)	843 (97%)	25 (3%)	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	188/201 (94%)	170 (90%)	18 (10%)	10	37
1	B	188/201 (94%)	178 (95%)	10 (5%)	26	65
1	E	188/201 (94%)	165 (88%)	23 (12%)	6	25
1	F	188/201 (94%)	174 (93%)	14 (7%)	16	52
All	All	752/804 (94%)	687 (91%)	65 (9%)	12	43

All (65) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	LEU
1	A	22	SER

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Mol	Chain	Res	Type
1	A	23	GLU
1	A	39	SER
1	A	56	ASP
1	A	89	SER
1	A	137	THR
1	A	140	LEU
1	A	146	THR
1	A	151	THR
1	A	159	SER
1	A	171	ARG
1	A	176	ASN
1	A	183	ASN
1	A	187	THR
1	A	192	VAL
1	A	208	THR
1	A	218	ASN
1	B	83	THR
1	B	118	ARG
1	B	119	HIS
1	B	120	ASN
1	B	133	ARG
1	B	171	ARG
1	B	175	TYR
1	B	195	HIS
1	B	200	LYS
1	B	207	ARG
1	E	44	HIS
1	E	54	LEU
1	E	56	ASP
1	E	60	LEU
1	E	81	ARG
1	E	82	ASP
1	E	85	GLU
1	E	100	VAL
1	E	105	LEU
1	E	117	ARG
1	E	128	SER
1	E	148	LEU
1	E	149	ASP
1	E	151	THR
1	E	171	ARG
1	E	179	TYR

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Mol	Chain	Res	Type
1	E	182	ASP
1	E	183	ASN
1	E	187	THR
1	E	194	ILE
1	E	203	LYS
1	E	212	PHE
1	E	216	LEU
1	F	22	SER
1	F	44	HIS
1	F	66	MET
1	F	71	MET
1	F	82	ASP
1	F	113	ARG
1	F	122	GLN
1	F	124	ASP
1	F	146	THR
1	F	153	LYS
1	F	183	ASN
1	F	184	GLU
1	F	197	LEU
1	F	210	ARG

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

Mol	Chain	Res	Type
1	A	119	HIS
1	E	120	ASN
1	E	183	ASN
1	F	14	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 carbohydrates in this entry.

5.6 Ligand geometry

Of 8 ligands modelled in this entry, 4 are monoatomic - leaving 4 for Mogul analysis.

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	BEF	A	301	1	0,3,3	0.00	-	0,3,3	0.00	-
4	BEF	B	301	1	0,3,3	0.00	-	0,3,3	0.00	-
4	BEF	E	301	1	0,3,3	0.00	-	0,3,3	0.00	-
4	BEF	F	301	1	0,3,3	0.00	-	0,3,3	0.00	-

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	BEF	A	301	1	-	0/0/0/0	0/0/0/0
4	BEF	B	301	1	-	0/0/0/0	0/0/0/0
4	BEF	E	301	1	-	0/0/0/0	0/0/0/0
4	BEF	F	301	1	-	0/0/0/0	0/0/0/0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	219/232 (94%)	-0.61	0 100 100	13, 37, 74, 111	0
1	B	219/232 (94%)	-0.07	11 (5%) 30 17	19, 73, 129, 144	0
1	E	219/232 (94%)	-0.41	3 (1%) 75 63	13, 45, 95, 154	0
1	F	219/232 (94%)	-0.34	6 (2%) 55 40	17, 48, 107, 146	0
2	C	25/25 (100%)	-0.28	0 100 100	30, 51, 101, 118	0
2	G	25/25 (100%)	-0.29	0 100 100	53, 64, 82, 138	0
3	D	25/25 (100%)	-0.29	0 100 100	37, 58, 100, 111	1 (4%)
3	H	25/25 (100%)	-0.38	0 100 100	42, 78, 93, 109	0
All	All	976/1028 (94%)	-0.35	20 (2%) 65 50	13, 49, 115, 154	1 (0%)

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	182	ASP	4.8
1	F	183	ASN	3.6
1	E	182	ASP	3.5
1	E	183	ASN	3.4
1	B	186	ALA	3.3
1	B	122	GLN	3.2
1	B	180	SER	3.0
1	B	123	GLY	2.9
1	B	145	GLU	2.8
1	B	219	ASN	2.7
1	F	121	ASN	2.6
1	F	182	ASP	2.5
1	B	183	ASN	2.5
1	B	137	THR	2.3
1	E	219	ASN	2.2
1	F	119	HIS	2.2

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Mol	Chain	Res	Type	RSRZ
1	F	219	ASN	2.2
1	B	120	ASN	2.0
1	B	124	ASP	2.0
1	F	120	ASN	2.0

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
4	BEF	E	301	4/4	0.77	0.30	2.53	52,52,52,52	0
4	BEF	A	301	4/4	0.86	0.24	1.87	60,60,61,61	0
4	BEF	F	301	4/4	0.89	0.20	0.92	58,58,58,59	0
4	BEF	B	301	4/4	0.93	0.19	0.69	55,56,56,57	0
5	MG	A	302	1/1	0.96	0.13	-0.60	29,29,29,29	0
5	MG	E	302	1/1	0.89	0.12	-0.97	12,12,12,12	0
5	MG	F	302	1/1	0.88	0.07	-1.81	33,33,33,33	0
5	MG	B	302	1/1	0.97	0.04	-2.25	40,40,40,40	0

6.5 Other polymers [i](#)

There are no such residues in this entry.