



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 15, 2017 – 08:01 pm GMT

PDB ID : 4APZ
Title : Structure of B. subtilis genomic dUTPase YncF in complex with dU, PPi and Mg in P1
Authors : Garcia-Nafria, J.; Timm, J.; Harrison, C.; Turkenburg, J.P.; Wilson, K.S.
Deposited on : 2012-04-11
Resolution : 2.01 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) : trunk28620

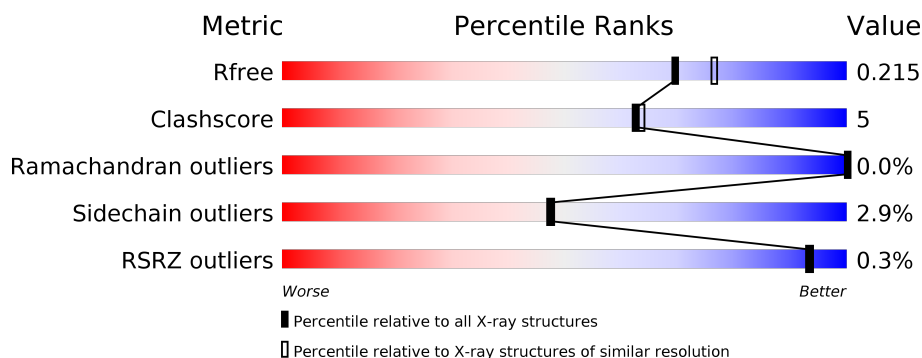
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.01 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	6609 (2.00-2.00)
Clashscore	112137	7775 (2.00-2.00)
Ramachandran outliers	110173	7679 (2.00-2.00)
Sidechain outliers	110143	7678 (2.00-2.00)
RSRZ outliers	101464	6696 (2.00-2.00)

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 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	1	144	<div> <div>%</div> <div> <div></div> <div>89%</div> <div>9%</div> <div>..</div> </div> </div>
1	2	144	<div> <div></div> <div>93%</div> <div>.</div> <div>..</div> </div>
1	3	144	<div> <div></div> <div>92%</div> <div>5%</div> <div>...</div> </div>
1	4	144	<div> <div></div> <div>89%</div> <div>8%</div> <div>..</div> </div>
1	5	144	<div> <div>%</div> <div> <div></div> <div>90%</div> <div>7%</div> <div>..</div> </div> </div>
1	6	144	<div> <div></div> <div>91%</div> <div>6%</div> <div>..</div> </div>





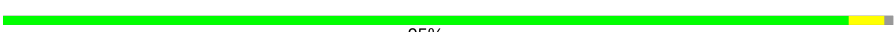









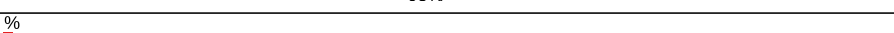
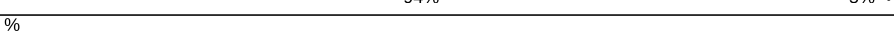
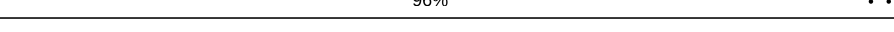
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Mol	Chain	Length	Quality of chain
1	7	144	 90% 8% ..
1	8	144	 84% 13% ..
1	9	144	 89% 8% ..
1	A	144	 89% 10% ..
1	B	144	 86% 12% ..
1	C	144	 88% 10% ..
1	D	144	 84% 14% ..
1	E	144	 85% 13% ..
1	F	144	 86% 12% ..
1	G	144	 86% 11% ..
1	H	144	 85% 11% ..
1	I	144	 87% 10% ..
1	J	144	 84% 14% ..
1	K	144	 77% 19% ..
1	L	144	 89% 10% .
1	M	144	 87% 10% ..
1	N	144	 90% 8% ..
1	O	144	 85% 13% ..
1	P	144	 88% 10% ..
1	Q	144	 84% 15% .
1	R	144	 87% 11% ..
1	S	144	 88% 8% ..
1	T	144	 90% 8% ..
1	U	144	 92% 6% ..
1	V	144	 88% 11% .

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Mol	Chain	Length	Quality of chain
1	W	144	 90% 9% ..
1	X	144	 92% 6% ..
1	Y	144	 89% 10% ..
1	Z	144	 87% 11% ..
1	a	144	 95% ..
1	b	144	 97% ..
1	c	144	 95% ..
1	d	144	 94% ..
1	e	144	 94% 5% ..
1	f	144	 94% 5% ..
1	g	144	 96% ..
1	h	144	 94% ..
1	i	144	 96% ..
1	j	144	 95% ..
1	k	144	 94% 5% ..
1	l	144	 96% ..
1	m	144	 95% ..

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
2	DUR	J	1145	X	-	-	-
5	PO4	l	1148	-	-	-	X

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 60911 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 PROBABLE DEOXYURIDINE 5'-TRIPHOSPHATE NUCLEOTIDOHYDROLASE YNCF.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	1	142	Total	C	N	O	S	9	2	0
			1148	730	193	218	7			
1	2	142	Total	C	N	O	S	0	2	0
			1149	730	194	218	7			
1	3	142	Total	C	N	O	S	3	0	0
			1136	721	191	218	6			
1	4	142	Total	C	N	O	S	0	4	0
			1162	739	196	220	7			
1	5	142	Total	C	N	O	S	7	1	0
			1141	725	191	218	7			
1	6	142	Total	C	N	O	S	11	1	0
			1141	725	191	218	7			
1	7	142	Total	C	N	O	S	0	0	0
			1136	721	191	218	6			
1	8	142	Total	C	N	O	S	0	1	0
			1141	725	191	218	7			
1	9	142	Total	C	N	O	S	3	2	0
			1150	731	193	219	7			
1	A	143	Total	C	N	O	S	4	2	0
			1151	731	192	221	7			
1	B	142	Total	C	N	O	S	4	2	0
			1148	730	193	218	7			
1	C	142	Total	C	N	O	S	6	2	0
			1148	729	193	220	6			
1	D	142	Total	C	N	O	S	7	1	0
			1141	725	191	218	7			
1	E	142	Total	C	N	O	S	3	1	0
			1143	726	193	218	6			
1	F	143	Total	C	N	O	S	3	2	0
			1155	734	194	220	7			
1	G	142	Total	C	N	O	S	6	2	0
			1152	732	195	219	6			

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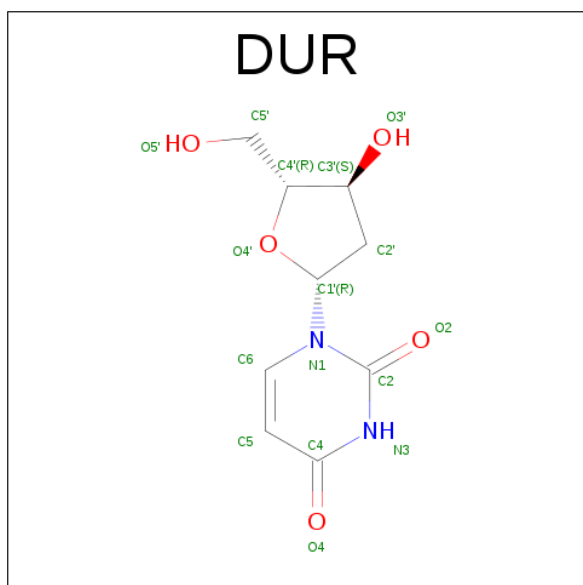
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	H	142	Total	C	N	O	S	3	0	0
			1136	721	191	218	6			
1	I	142	Total	C	N	O	S	7	1	0
			1141	725	191	218	7			
1	J	142	Total	C	N	O	S	2	1	0
			1143	726	193	218	6			
1	K	142	Total	C	N	O	S	11	1	0
			1142	725	191	220	6			
1	L	142	Total	C	N	O	S	0	0	0
			1136	721	191	218	6			
1	M	142	Total	C	N	O	S	2	1	0
			1139	723	191	219	6			
1	N	142	Total	C	N	O	S	3	0	0
			1136	721	191	218	6			
1	O	142	Total	C	N	O	S	0	0	0
			1136	721	191	218	6			
1	P	142	Total	C	N	O	S	2	2	0
			1150	731	193	219	7			
1	Q	143	Total	C	N	O	S	0	1	0
			1150	730	194	220	6			
1	R	142	Total	C	N	O	S	3	1	0
			1143	726	193	218	6			
1	S	142	Total	C	N	O	S	0	2	0
			1148	730	193	218	7			
1	T	142	Total	C	N	O	S	4	2	0
			1148	730	193	218	7			
1	U	142	Total	C	N	O	S	0	2	0
			1148	730	193	218	7			
1	V	142	Total	C	N	O	S	3	1	0
			1141	725	191	218	7			
1	W	142	Total	C	N	O	S	0	0	0
			1136	721	191	218	6			
1	X	142	Total	C	N	O	S	0	3	0
			1151	732	193	219	7			
1	Y	142	Total	C	N	O	S	3	1	0
			1141	725	191	218	7			
1	Z	142	Total	C	N	O	S	8	2	0
			1150	731	193	219	7			
1	a	143	Total	C	N	O	S	7	1	0
			1148	729	192	220	7			
1	b	143	Total	C	N	O	S	3	0	0
			1143	725	192	220	6			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	c	142	Total	C	N	O	S	8	2	0
			1148	730	193	218	7			
1	d	142	Total	C	N	O	S	0	1	0
			1141	725	191	218	7			
1	e	142	Total	C	N	O	S	0	1	0
			1141	725	191	218	7			
1	f	142	Total	C	N	O	S	0	1	0
			1141	725	191	218	7			
1	g	142	Total	C	N	O	S	3	2	0
			1153	732	194	220	7			
1	h	142	Total	C	N	O	S	4	0	0
			1136	721	191	218	6			
1	i	142	Total	C	N	O	S	4	1	0
			1141	725	191	218	7			
1	j	143	Total	C	N	O	S	2	1	0
			1148	729	192	220	7			
1	k	142	Total	C	N	O	S	4	1	0
			1141	725	191	218	7			
1	l	142	Total	C	N	O	S	0	2	0
			1148	730	193	218	7			
1	m	142	Total	C	N	O	S	0	1	0
			1141	725	191	218	7			

- Molecule 2 is 2'-DEOXYURIDINE (three-letter code: DUR) (formula: C₉H₁₂N₂O₅).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	1	1	Total	C	N	O	0	0
			16	9	2	5		
2	2	1	Total	C	N	O	0	0
			16	9	2	5		
2	3	1	Total	C	N	O	0	0
			16	9	2	5		
2	4	1	Total	C	N	O	0	0
			16	9	2	5		
2	5	1	Total	C	N	O	0	0
			16	9	2	5		
2	6	1	Total	C	N	O	0	0
			16	9	2	5		
2	7	1	Total	C	N	O	0	0
			16	9	2	5		
2	8	1	Total	C	N	O	0	0
			16	9	2	5		
2	9	1	Total	C	N	O	0	0
			16	9	2	5		
2	A	1	Total	C	N	O	0	0
			16	9	2	5		
2	B	1	Total	C	N	O	0	0
			16	9	2	5		
2	C	1	Total	C	N	O	0	0
			16	9	2	5		
2	D	1	Total	C	N	O	0	0
			16	9	2	5		
2	E	1	Total	C	N	O	0	0
			16	9	2	5		
2	F	1	Total	C	N	O	0	0
			16	9	2	5		
2	G	1	Total	C	N	O	0	0
			16	9	2	5		
2	H	1	Total	C	N	O	0	0
			16	9	2	5		
2	I	1	Total	C	N	O	0	0
			16	9	2	5		
2	J	1	Total	C	N	O	0	0
			16	9	2	5		
2	K	1	Total	C	N	O	0	0
			16	9	2	5		
2	L	1	Total	C	N	O	0	0
			16	9	2	5		
2	M	1	Total	C	N	O	0	0
			16	9	2	5		

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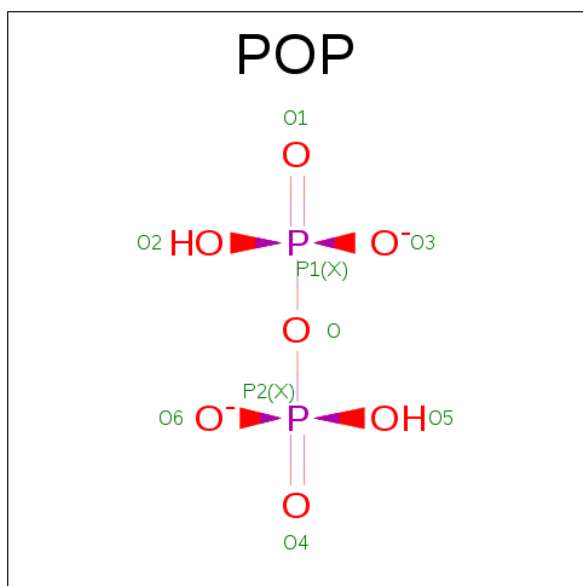
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	N	1	Total	C	N	O	0	0
			16	9	2	5		
2	O	1	Total	C	N	O	0	0
			16	9	2	5		
2	P	1	Total	C	N	O	0	0
			16	9	2	5		
2	Q	1	Total	C	N	O	0	0
			16	9	2	5		
2	R	1	Total	C	N	O	0	0
			16	9	2	5		
2	S	1	Total	C	N	O	0	0
			16	9	2	5		
2	T	1	Total	C	N	O	0	0
			16	9	2	5		
2	U	1	Total	C	N	O	0	0
			16	9	2	5		
2	V	1	Total	C	N	O	0	0
			16	9	2	5		
2	W	1	Total	C	N	O	0	0
			16	9	2	5		
2	X	1	Total	C	N	O	0	0
			16	9	2	5		
2	Y	1	Total	C	N	O	0	0
			16	9	2	5		
2	Z	1	Total	C	N	O	0	0
			16	9	2	5		
2	a	1	Total	C	N	O	0	0
			16	9	2	5		
2	b	1	Total	C	N	O	0	0
			16	9	2	5		
2	c	1	Total	C	N	O	0	0
			16	9	2	5		
2	d	1	Total	C	N	O	0	0
			16	9	2	5		
2	e	1	Total	C	N	O	0	0
			16	9	2	5		
2	f	1	Total	C	N	O	0	0
			16	9	2	5		
2	g	1	Total	C	N	O	0	0
			16	9	2	5		
2	h	1	Total	C	N	O	0	0
			16	9	2	5		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	i	1	Total	C	N	O	0	0
			16	9	2	5		
2	j	1	Total	C	N	O	0	0
			16	9	2	5		
2	k	1	Total	C	N	O	0	0
			16	9	2	5		
2	l	1	Total	C	N	O	0	0
			16	9	2	5		
2	m	1	Total	C	N	O	0	0
			16	9	2	5		

- Molecule 3 is PYROPHOSPHATE 2- (three-letter code: POP) (formula: $\text{H}_2\text{O}_7\text{P}_2$).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	1	1	Total	O P	0	0
			9 7 2			
3	2	1	Total	O P	0	0
			9 7 2			
3	3	1	Total	O P	0	0
			9 7 2			
3	4	1	Total	O P	0	0
			9 7 2			
3	5	1	Total	O P	0	0
			9 7 2			
3	6	1	Total	O P	0	0
			9 7 2			

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	7	1	Total 9	O 7	P 2	0	0
3	8	1	Total 9	O 7	P 2	0	0
3	9	1	Total 9	O 7	P 2	0	0
3	A	1	Total 9	O 7	P 2	0	0
3	B	1	Total 9	O 7	P 2	0	0
3	C	1	Total 9	O 7	P 2	0	0
3	D	1	Total 9	O 7	P 2	0	0
3	E	1	Total 9	O 7	P 2	0	0
3	F	1	Total 9	O 7	P 2	0	0
3	G	1	Total 9	O 7	P 2	0	0
3	H	1	Total 9	O 7	P 2	0	0
3	I	1	Total 9	O 7	P 2	0	0
3	J	1	Total 9	O 7	P 2	0	0
3	K	1	Total 9	O 7	P 2	0	0
3	L	1	Total 9	O 7	P 2	0	0
3	M	1	Total 9	O 7	P 2	0	0
3	N	1	Total 9	O 7	P 2	0	0
3	O	1	Total 9	O 7	P 2	0	0
3	P	1	Total 9	O 7	P 2	0	0
3	Q	1	Total 9	O 7	P 2	0	0
3	R	1	Total 9	O 7	P 2	0	0

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	S	1	Total	O	P	0	0
			9	7	2		
3	T	1	Total	O	P	0	0
			9	7	2		
3	U	1	Total	O	P	0	0
			9	7	2		
3	V	1	Total	O	P	0	0
			9	7	2		
3	W	1	Total	O	P	0	0
			9	7	2		
3	X	1	Total	O	P	0	0
			9	7	2		
3	Y	1	Total	O	P	0	0
			9	7	2		
3	Z	1	Total	O	P	0	0
			9	7	2		
3	a	1	Total	O	P	0	0
			9	7	2		
3	b	1	Total	O	P	0	0
			9	7	2		
3	c	1	Total	O	P	0	0
			9	7	2		
3	d	1	Total	O	P	0	0
			9	7	2		
3	e	1	Total	O	P	0	0
			9	7	2		
3	f	1	Total	O	P	0	0
			9	7	2		
3	g	1	Total	O	P	0	0
			9	7	2		
3	h	1	Total	O	P	0	0
			9	7	2		
3	i	1	Total	O	P	0	0
			9	7	2		
3	j	1	Total	O	P	0	0
			9	7	2		
3	k	1	Total	O	P	0	0
			9	7	2		
3	l	1	Total	O	P	0	0
			9	7	2		
3	m	1	Total	O	P	0	0
			9	7	2		

- Molecule 4 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	P	1	Total 1	Mg 1	0	0
4	g	1	Total 1	Mg 1	0	0
4	K	1	Total 1	Mg 1	0	0
4	h	1	Total 1	Mg 1	0	0
4	B	1	Total 1	Mg 1	0	0
4	c	1	Total 1	Mg 1	0	0
4	6	1	Total 1	Mg 1	0	0
4	W	1	Total 1	Mg 1	0	0
4	N	1	Total 1	Mg 1	0	0
4	X	1	Total 1	Mg 1	0	0
4	2	1	Total 1	Mg 1	0	0
4	S	1	Total 1	Mg 1	0	0
4	f	1	Total 1	Mg 1	0	0
4	J	1	Total 1	Mg 1	0	0
4	k	1	Total 1	Mg 1	0	0
4	E	1	Total 1	Mg 1	0	0
4	b	1	Total 1	Mg 1	0	0
4	V	1	Total 1	Mg 1	0	0
4	A	1	Total 1	Mg 1	0	0
4	5	1	Total 1	Mg 1	0	0
4	R	1	Total 1	Mg 1	0	0

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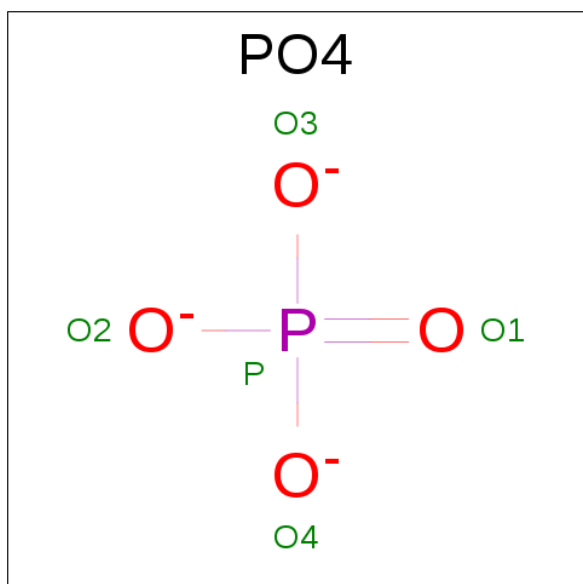
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	M	1	Total 1	Mg 1	0	0
4	j	1	Total 1	Mg 1	0	0
4	1	1	Total 1	Mg 1	0	0
4	D	1	Total 1	Mg 1	0	0
4	e	1	Total 1	Mg 1	0	0
4	I	1	Total 1	Mg 1	0	0
4	Z	1	Total 1	Mg 1	0	0
4	a	1	Total 1	Mg 1	0	0
4	4	1	Total 1	Mg 1	0	0
4	U	1	Total 1	Mg 1	0	0
4	9	1	Total 1	Mg 1	0	0
4	L	1	Total 1	Mg 1	0	0
4	m	1	Total 1	Mg 1	0	0
4	G	1	Total 1	Mg 1	0	0
4	Q	1	Total 1	Mg 1	0	0
4	d	1	Total 1	Mg 1	0	0
4	H	1	Total 1	Mg 1	0	0
4	i	1	Total 1	Mg 1	0	0
4	C	1	Total 1	Mg 1	0	0
4	7	1	Total 1	Mg 1	0	0
4	T	1	Total 1	Mg 1	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	8	1	Total	Mg	0	0
			1	1		
4	O	1	Total	Mg	0	0
			1	1		
4	Y	1	Total	Mg	0	0
			1	1		
4	l	1	Total	Mg	0	0
			1	1		
4	3	1	Total	Mg	0	0
			1	1		
4	F	1	Total	Mg	0	0
			1	1		

- Molecule 5 is PHOSPHATE ION (three-letter code: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	e	1	Total	O	P	0	0
			5	4	1		
5	l	1	Total	O	P	0	0
			5	4	1		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	1	120	Total	O	0	0
			120	120		

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	2	139	Total O 139 139	0	0
6	3	102	Total O 102 102	0	0
6	4	85	Total O 85 85	0	0
6	5	109	Total O 109 109	0	0
6	6	93	Total O 93 93	0	0
6	7	69	Total O 69 69	0	0
6	8	136	Total O 136 136	0	0
6	9	101	Total O 101 101	0	0
6	A	122	Total O 122 122	0	0
6	a	1	Total O 1 1	0	0
6	A	3	Total O 3 3	0	0
6	B	96	Total O 96 96	0	0
6	C	67	Total O 67 67	0	0
6	D	133	Total O 133 133	0	0
6	E	111	Total O 111 111	0	0
6	F	86	Total O 86 86	0	0
6	G	141	Total O 141 141	0	0
6	H	123	Total O 123 123	0	0
6	I	84	Total O 84 84	0	0
6	J	116	Total O 116 116	0	0
6	K	95	Total O 95 95	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	L	70	Total 70	O 70	0	0
6	M	124	Total 124	O 124	0	0
6	N	98	Total 98	O 98	0	0
6	O	68	Total 68	O 68	0	0
6	P	142	Total 142	O 142	0	0
6	Q	103	Total 103	O 103	0	0
6	R	75	Total 75	O 75	0	0
6	S	133	Total 133	O 133	0	0
6	T	90	Total 90	O 90	0	0
6	U	62	Total 62	O 62	0	0
6	V	118	Total 118	O 118	0	0
6	W	97	Total 97	O 97	0	0
6	X	70	Total 70	O 70	0	0
6	Y	94	Total 94	O 94	0	0
6	Z	60	Total 60	O 60	0	0
6	a	69	Total 69	O 69	0	0
6	b	127	Total 127	O 127	0	0
6	c	105	Total 105	O 105	0	0
6	d	63	Total 63	O 63	0	0
6	e	128	Total 128	O 128	0	0
6	f	92	Total 92	O 92	0	0

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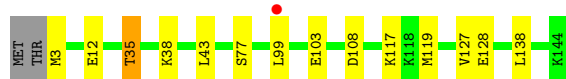
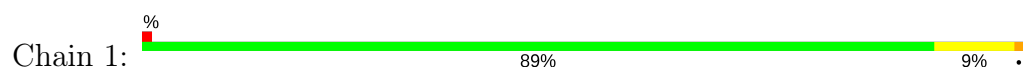
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	g	74	Total 74	O 74	0	0
6	h	143	Total 143	O 143	0	0
6	i	99	Total 99	O 99	0	0
6	j	63	Total 63	O 63	0	0
6	k	109	Total 109	O 109	0	0
6	l	69	Total 69	O 69	0	0
6	m	49	Total 49	O 49	0	0

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.

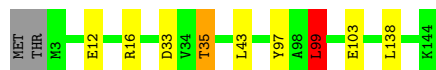
- Molecule 1: PROBABLE DEOXYURIDINE 5'-TRIPHOSPHATE NUCLEOTIDOHYDROLASE YNCF



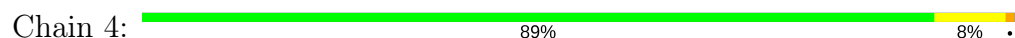
- Molecule 1: PROBABLE DEOXYURIDINE 5'-TRIPHOSPHATE NUCLEOTIDOHYDROLASE YNCF



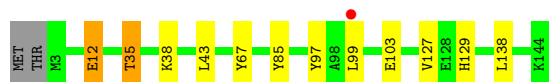
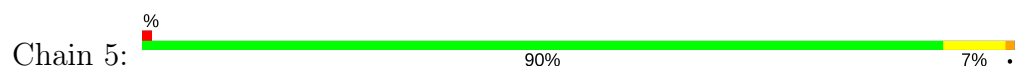
- Molecule 1: PROBABLE DEOXYURIDINE 5'-TRIPHOSPHATE NUCLEOTIDOHYDROLASE YNCF



- Molecule 1: PROBABLE DEOXYURIDINE 5'-TRIPHOSPHATE NUCLEOTIDOHYDROLASE YNCF

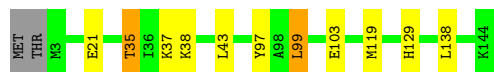


- Molecule 1: PROBABLE DEOXYURIDINE 5'-TRIPHOSPHATE NUCLEOTIDOHYDROLASE YNCF



- Molecule 1: PROBABLE DEOXYURIDINE 5'-TRIPHOSPHATE NUCLEOTIDOHYDROLASE YNCF

Chain 6:  91% 6% ..




- Molecule 1: PROBABLE DEOXYURIDINE 5'-TRIPHOSPHATE NUCLEOTIDOHYDROLASE YNCF

Chain 7:  90% 8% ..



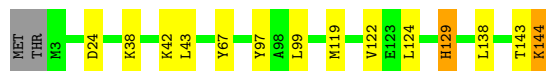
- Molecule 1: PROBABLE DEOXYURIDINE 5'-TRIPHOSPHATE NUCLEOTIDOHYDROLASE YNCF

Chain 8:  84% 13% ..




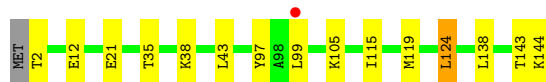
- Molecule 1: PROBABLE DEOXYURIDINE 5'-TRIPHOSPHATE NUCLEOTIDOHYDROLASE YNCF

Chain 9:  89% 8% ..



- Molecule 1: PROBABLE DEOXYURIDINE 5'-TRIPHOSPHATE NUCLEOTIDOHYDROLASE YNCF

Chain A:  89% 10% ..

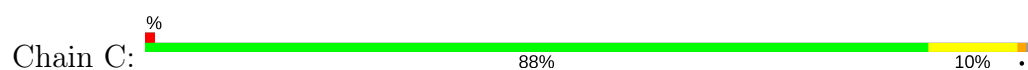


- Molecule 1: PROBABLE DEOXYURIDINE 5'-TRIPHOSPHATE NUCLEOTIDOHYDROLASE YNCF

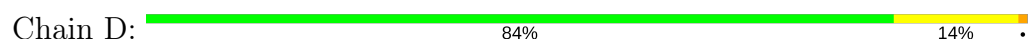
Chain B:  86% 12% ..



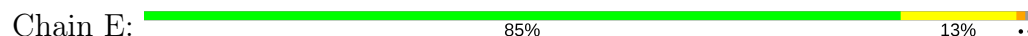
- Molecule 1: PROBABLE DEOXYURIDINE 5'-TRIPHOSPHATE NUCLEOTIDOHYDROLASE YNCF



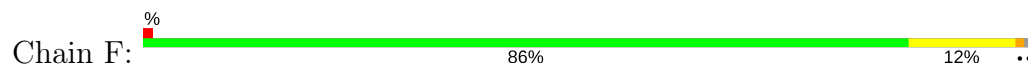
- Molecule 1: PROBABLE DEOXYURIDINE 5'-TRIPHOSPHATE NUCLEOTIDOHYDROLASE YNCF



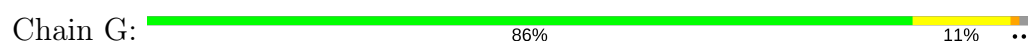
- Molecule 1: PROBABLE DEOXYURIDINE 5'-TRIPHOSPHATE NUCLEOTIDOHYDROLASE YNCF



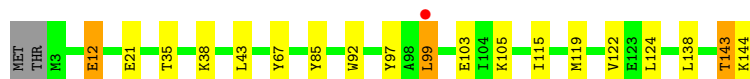
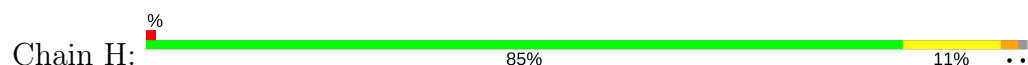
- Molecule 1: PROBABLE DEOXYURIDINE 5'-TRIPHOSPHATE NUCLEOTIDOHYDROLASE YNCF



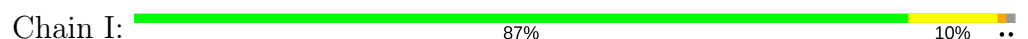
- Molecule 1: PROBABLE DEOXYURIDINE 5'-TRIPHOSPHATE NUCLEOTIDOHYDROLASE YNCF

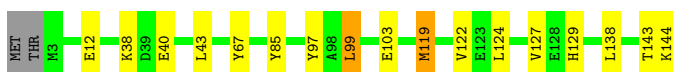


- Molecule 1: PROBABLE DEOXYURIDINE 5'-TRIPHOSPHATE NUCLEOTIDOHYDROLASE YNCF



- Molecule 1: PROBABLE DEOXYURIDINE 5'-TRIPHOSPHATE NUCLEOTIDOHYDROLASE YNCF





- Molecule 1: PROBABLE DEOXYURIDINE 5'-TRIPHOSPHATE NUCLEOTIDOHYDROLASE YNCF

Chain J: 84% 14% ..



- Molecule 1: PROBABLE DEOXYURIDINE 5'-TRIPHOSPHATE NUCLEOTIDOHYDROLASE YNCF

Chain K: 77% 19% ..



- Molecule 1: PROBABLE DEOXYURIDINE 5'-TRIPHOSPHATE NUCLEOTIDOHYDROLASE YNCF

Chain L: 89% 10% .



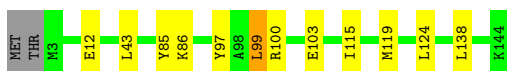
- Molecule 1: PROBABLE DEOXYURIDINE 5'-TRIPHOSPHATE NUCLEOTIDOHYDROLASE YNCF

Chain M: 87% 10% ..



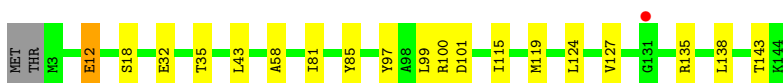
- Molecule 1: PROBABLE DEOXYURIDINE 5'-TRIPHOSPHATE NUCLEOTIDOHYDROLASE YNCF

Chain N: 90% 8% ..




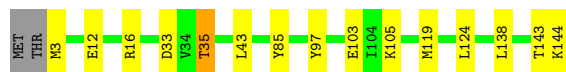
- Molecule 1: PROBABLE DEOXYURIDINE 5'-TRIPHOSPHATE NUCLEOTIDOHYDROLASE YNCF

Chain O: 85% 13% ..




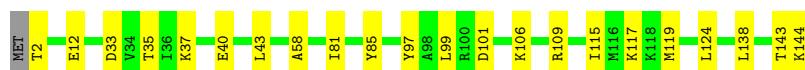
- Molecule 1: PROBABLE DEOXYURIDINE 5'-TRIPHOSPHATE NUCLEOTIDOHYDROLASE YNCF

Chain P:  88% 10% ..



- Molecule 1: PROBABLE DEOXYURIDINE 5'-TRIPHOSPHATE NUCLEOTIDOHYDROLASE YNCF

Chain Q:  84% 15% .




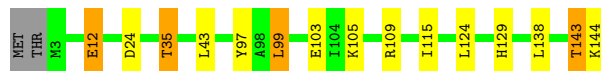
- Molecule 1: PROBABLE DEOXYURIDINE 5'-TRIPHOSPHATE NUCLEOTIDOHYDROLASE YNCF

Chain R:  87% 11% ..



- Molecule 1: PROBABLE DEOXYURIDINE 5'-TRIPHOSPHATE NUCLEOTIDOHYDROLASE YNCF

Chain S:  88% 8% ..



- Molecule 1: PROBABLE DEOXYURIDINE 5'-TRIPHOSPHATE NUCLEOTIDOHYDROLASE YNCF

Chain T:  90% 8% ..




- Molecule 1: PROBABLE DEOXYURIDINE 5'-TRIPHOSPHATE NUCLEOTIDOHYDROLASE YNCF

Chain U:  92% 6% ..



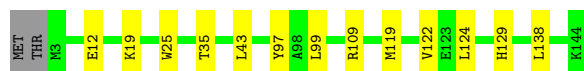
- Molecule 1: PROBABLE DEOXYURIDINE 5'-TRIPHOSPHATE NUCLEOTIDOHYDROLASE YNCF

Chain V:  88% 11%



- Molecule 1: PROBABLE DEOXYURIDINE 5'-TRIPHOSPHATE NUCLEOTIDOHYDROLASE YNCF

Chain W:  90% 9%




- Molecule 1: PROBABLE DEOXYURIDINE 5'-TRIPHOSPHATE NUCLEOTIDOHYDROLASE YNCF

Chain X:  92% 6%




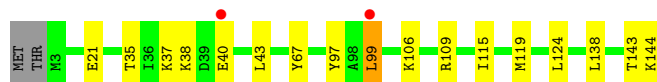
- Molecule 1: PROBABLE DEOXYURIDINE 5'-TRIPHOSPHATE NUCLEOTIDOHYDROLASE YNCF

Chain Y:  89% 10%



- Molecule 1: PROBABLE DEOXYURIDINE 5'-TRIPHOSPHATE NUCLEOTIDOHYDROLASE YNCF

Chain Z:  87% 11%



- Molecule 1: PROBABLE DEOXYURIDINE 5'-TRIPHOSPHATE NUCLEOTIDOHYDROLASE YNCF

Chain a:  95%



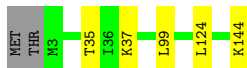
- Molecule 1: PROBABLE DEOXYURIDINE 5'-TRIPHOSPHATE NUCLEOTIDOHYDROLASE YNCF

Chain b:  97%



- Molecule 1: PROBABLE DEOXYURIDINE 5'-TRIPHOSPHATE NUCLEOTIDOHYDROLASE YNCF

Chain c:  95% ..



- Molecule 1: PROBABLE DEOXYURIDINE 5'-TRIPHOSPHATE NUCLEOTIDOHYDROLASE YNCF

Chain d:  94% ..



- Molecule 1: PROBABLE DEOXYURIDINE 5'-TRIPHOSPHATE NUCLEOTIDOHYDROLASE YNCF

Chain e:  94% 5% .



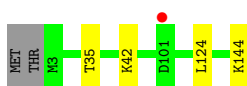
- Molecule 1: PROBABLE DEOXYURIDINE 5'-TRIPHOSPHATE NUCLEOTIDOHYDROLASE YNCF

Chain f:  94% 5% .



- Molecule 1: PROBABLE DEOXYURIDINE 5'-TRIPHOSPHATE NUCLEOTIDOHYDROLASE YNCF

Chain g:  96% ..



- Molecule 1: PROBABLE DEOXYURIDINE 5'-TRIPHOSPHATE NUCLEOTIDOHYDROLASE YNCF

Chain h:  94% ..



- Molecule 1: PROBABLE DEOXYURIDINE 5'-TRIPHOSPHATE NUCLEOTIDOHYDROLASE YNCF

Chain i:  96%



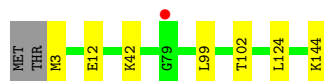
- Molecule 1: PROBABLE DEOXYURIDINE 5'-TRIPHOSPHATE NUCLEOTIDOHYDROLASE YNCF

Chain j:  95%



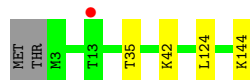
- Molecule 1: PROBABLE DEOXYURIDINE 5'-TRIPHOSPHATE NUCLEOTIDOHYDROLASE YNCF

Chain k:  94%



- Molecule 1: PROBABLE DEOXYURIDINE 5'-TRIPHOSPHATE NUCLEOTIDOHYDROLASE YNCF

Chain l:  96%



- Molecule 1: PROBABLE DEOXYURIDINE 5'-TRIPHOSPHATE NUCLEOTIDOHYDROLASE YNCF

Chain m:  95%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	97.06Å 97.07Å 193.85Å 89.72° 88.47° 90.11°	Depositor
Resolution (Å)	36.44 – 2.01 36.44 – 2.01	Depositor EDS
% Data completeness (in resolution range)	96.7 (36.44-2.01) 96.6 (36.44-2.01)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.29 (at 2.01Å)	Xtriage
Refinement program	REFMAC 5.6.0119	Depositor
R, R_{free}	0.169 , 0.212 0.174 , 0.215	Depositor DCC
R_{free} test set	22882 reflections (5.29%)	DCC
Wilson B-factor (Å ²)	23.3	Xtriage
Anisotropy	0.284	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 48.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.43$, $\langle L^2 \rangle = 0.26$	Xtriage
Estimated twinning fraction	0.044 for k,-h,l 0.044 for -k,h,l 0.046 for h,-k,-l 0.048 for -h,k,-l 0.044 for -h,-k,l 0.039 for k,h,-l 0.115 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	60911	wwPDB-VP
Average B, all atoms (Å ²)	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.99% 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, DUR, PO4, POP

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	1	0.81	0/1176	0.82	1/1582 (0.1%)
1	2	0.88	0/1176	0.85	0/1581
1	3	0.94	0/1157	0.89	2/1556 (0.1%)
1	4	0.96	1/1196 (0.1%)	0.90	3/1608 (0.2%)
1	5	0.85	0/1165	0.85	1/1567 (0.1%)
1	6	0.87	0/1165	0.86	1/1567 (0.1%)
1	7	0.82	0/1157	0.83	1/1556 (0.1%)
1	8	0.88	0/1165	0.86	2/1567 (0.1%)
1	9	0.90	0/1173	0.85	1/1575 (0.1%)
1	A	0.83	0/1178	0.82	1/1585 (0.1%)
1	B	0.86	0/1176	0.84	0/1582
1	C	0.80	0/1171	0.82	0/1572
1	D	0.90	1/1165 (0.1%)	0.88	1/1567 (0.1%)
1	E	0.90	0/1168	0.91	2/1571 (0.1%)
1	F	0.92	1/1183 (0.1%)	0.88	0/1592
1	G	0.92	0/1176	0.89	1/1579 (0.1%)
1	H	0.99	1/1157 (0.1%)	0.91	1/1556 (0.1%)
1	I	0.94	2/1165 (0.2%)	0.92	1/1567 (0.1%)
1	J	0.90	0/1168	0.92	5/1571 (0.3%)
1	K	0.86	0/1166	0.88	1/1568 (0.1%)
1	L	0.89	1/1157 (0.1%)	0.87	0/1556
1	M	0.90	0/1163	0.85	1/1564 (0.1%)
1	N	0.83	0/1157	0.86	2/1556 (0.1%)
1	O	0.90	0/1157	0.85	1/1556 (0.1%)
1	P	0.98	0/1173	0.88	1/1575 (0.1%)
1	Q	0.93	0/1175	0.91	2/1581 (0.1%)
1	R	0.94	0/1168	0.87	2/1571 (0.1%)
1	S	0.87	0/1176	0.86	3/1582 (0.2%)
1	T	0.86	0/1176	0.85	1/1582 (0.1%)
1	U	0.87	0/1176	0.86	0/1582
1	V	0.96	2/1165 (0.2%)	0.91	0/1567
1	W	0.92	2/1157 (0.2%)	0.85	2/1556 (0.1%)

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	X	0.90	0/1182	0.91	1/1590 (0.1%)
1	Y	0.87	0/1165	0.88	0/1567
1	Z	0.85	0/1173	0.84	1/1575 (0.1%)
1	a	0.88	0/1172	0.89	2/1577 (0.1%)
1	b	0.87	0/1164	0.86	0/1566
1	c	0.89	1/1176 (0.1%)	0.87	1/1582 (0.1%)
1	d	0.89	1/1165 (0.1%)	0.88	1/1567 (0.1%)
1	e	0.92	2/1165 (0.2%)	0.86	0/1567
1	f	0.93	1/1165 (0.1%)	0.89	2/1567 (0.1%)
1	g	0.89	0/1173	0.85	0/1575
1	h	0.94	1/1157 (0.1%)	0.88	1/1556 (0.1%)
1	i	0.94	1/1165 (0.1%)	0.87	0/1567
1	j	0.93	0/1172	0.88	0/1577
1	k	0.92	0/1165	0.99	0/1567
1	l	0.81	0/1176	0.82	0/1582
1	m	0.81	1/1165 (0.1%)	0.85	1/1567 (0.1%)
All	All	0.89	19/56103 (0.0%)	0.87	50/75446 (0.1%)

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	W	25	TRP	CD2-CE2	5.90	1.48	1.41
1	m	25	TRP	CD2-CE2	5.78	1.48	1.41
1	F	25	TRP	CD2-CE2	5.77	1.48	1.41
1	c	37	LYS	CB-CG	-5.76	1.36	1.52
1	I	12	GLU	CG-CD	5.75	1.60	1.51
1	f	92	TRP	CD2-CE2	5.43	1.47	1.41
1	i	92	TRP	CD2-CE2	5.34	1.47	1.41
1	L	128	GLU	CG-CD	5.31	1.59	1.51
1	e	92	TRP	CD2-CE2	5.27	1.47	1.41
1	H	92	TRP	CD2-CE2	5.25	1.47	1.41
1	h	54	GLU	CB-CG	-5.23	1.42	1.52
1	d	25	TRP	CD2-CE2	5.22	1.47	1.41
1	V	25	TRP	NE1-CE2	-5.17	1.30	1.37
1	4	25	TRP	CD2-CE2	5.12	1.47	1.41
1	D	18	SER	CB-OG	5.08	1.48	1.42
1	W	25	TRP	NE1-CE2	-5.07	1.30	1.37
1	I	40	GLU	CG-CD	5.06	1.59	1.51
1	e	40	GLU	CD-OE1	5.05	1.31	1.25
1	V	92	TRP	CD2-CE2	5.01	1.47	1.41

All (50) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	J	134	ASP	CB-CG-OD2	-6.00	112.90	118.30
1	K	109	ARG	NE-CZ-NH1	5.93	123.26	120.30
1	J	100	ARG	NE-CZ-NH2	-5.89	117.36	120.30
1	f	40	GLU	CA-CB-CG	5.88	126.33	113.40
1	J	134	ASP	CB-CG-OD1	5.87	123.58	118.30
1	3	99	LEU	CB-CG-CD1	5.86	120.96	111.00
1	W	109	ARG	NE-CZ-NH2	-5.80	117.40	120.30
1	4	109	ARG	NE-CZ-NH1	5.78	123.19	120.30
1	N	99	LEU	CB-CG-CD1	5.76	120.80	111.00
1	E	29	ARG	NE-CZ-NH1	5.73	123.16	120.30
1	E	135	ARG	NE-CZ-NH1	5.71	123.15	120.30
1	S	109	ARG	NE-CZ-NH1	5.68	123.14	120.30
1	P	33	ASP	CB-CG-OD1	5.64	123.37	118.30
1	a	99	LEU	CB-CG-CD1	5.62	120.55	111.00
1	4	6	LYS	CD-CE-NZ	-5.61	98.80	111.70
1	G	109	ARG	NE-CZ-NH1	5.61	123.11	120.30
1	d	99	LEU	CB-CG-CD2	5.61	120.54	111.00
1	I	99	LEU	CB-CG-CD1	5.55	120.44	111.00
1	R	109	ARG	NE-CZ-NH2	-5.54	117.53	120.30
1	S	24	ASP	CB-CG-OD1	5.54	123.28	118.30
1	6	99	LEU	CB-CG-CD2	5.53	120.40	111.00
1	4	6	LYS	CG-CD-CE	5.50	128.39	111.90
1	Q	33	ASP	CB-CG-OD1	5.48	123.23	118.30
1	9	24	ASP	CB-CG-OD1	5.47	123.22	118.30
1	O	135	ARG	NE-CZ-NH1	5.38	122.99	120.30
1	a	33	ASP	CB-CG-OD1	5.36	123.12	118.30
1	Q	109	ARG	NE-CZ-NH2	-5.35	117.62	120.30
1	M	100	ARG	NE-CZ-NH2	-5.31	117.65	120.30
1	3	33	ASP	CB-CG-OD1	5.29	123.06	118.30
1	H	105	LYS	CD-CE-NZ	-5.29	99.52	111.70
1	5	12	GLU	CA-CB-CG	5.29	125.03	113.40
1	f	99	LEU	CB-CG-CD1	5.27	119.97	111.00
1	S	99	LEU	CB-CG-CD1	5.26	119.94	111.00
1	A	124	LEU	CB-CG-CD1	5.25	119.92	111.00
1	8	109	ARG	NE-CZ-NH1	5.24	122.92	120.30
1	Z	109	ARG	NE-CZ-NH1	5.23	122.91	120.30
1	h	119	MET	CG-SD-CE	5.21	108.54	100.20
1	D	78	MET	CG-SD-CE	-5.21	91.87	100.20
1	R	100	ARG	NE-CZ-NH2	-5.19	117.70	120.30
1	1	108	ASP	CB-CG-OD2	-5.16	113.66	118.30
1	J	119	MET	CG-SD-CE	-5.16	91.95	100.20
1	J	33	ASP	CB-CG-OD1	5.16	122.94	118.30
1	c	99	LEU	CB-CG-CD1	5.14	119.73	111.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	N	100	ARG	NE-CZ-NH1	5.04	122.82	120.30
1	T	99	LEU	CB-CG-CD1	5.04	119.57	111.00
1	7	134	ASP	CB-CG-OD1	5.04	122.84	118.30
1	W	99	LEU	CB-CG-CD1	5.03	119.55	111.00
1	m	99	LEU	CB-CG-CD1	5.03	119.56	111.00
1	X	99	LEU	CB-CG-CD1	5.02	119.54	111.00
1	8	101	ASP	CB-CG-OD2	-5.02	113.78	118.30

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	1	1148	0	1146	22	0
1	2	1149	0	1152	15	0
1	3	1136	0	1130	7	0
1	4	1162	0	1165	12	0
1	5	1141	0	1139	21	0
1	6	1141	0	1139	15	0
1	7	1136	0	1130	20	0
1	8	1141	0	1139	23	0
1	9	1150	0	1152	18	0
1	A	1151	0	1151	36	0
1	B	1148	0	1146	31	0
1	C	1148	0	1148	28	0
1	D	1141	0	1139	27	0
1	E	1143	0	1137	28	0
1	F	1155	0	1153	35	0
1	G	1152	0	1150	25	0
1	H	1136	0	1130	24	0
1	I	1141	0	1139	23	0
1	J	1143	0	1137	24	0
1	K	1142	0	1136	29	0
1	L	1136	0	1130	14	0
1	M	1139	0	1135	26	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	N	1136	0	1130	16	0
1	O	1136	0	1130	21	0
1	P	1150	0	1152	19	0
1	Q	1150	0	1144	22	0
1	R	1143	0	1137	19	0
1	S	1148	0	1146	14	0
1	T	1148	0	1146	11	0
1	U	1148	0	1146	11	0
1	V	1141	0	1139	11	0
1	W	1136	0	1130	9	0
1	X	1151	0	1151	9	0
1	Y	1141	0	1139	19	0
1	Z	1150	0	1152	21	0
1	a	1148	0	1146	0	0
1	b	1143	0	1137	0	0
1	c	1148	0	1146	0	0
1	d	1141	0	1139	0	0
1	e	1141	0	1139	0	0
1	f	1141	0	1139	0	0
1	g	1153	0	1152	0	0
1	h	1136	0	1130	0	0
1	i	1141	0	1139	0	0
1	j	1148	0	1146	0	0
1	k	1141	0	1139	0	0
1	l	1148	0	1146	0	0
1	m	1141	0	1139	0	0
2	1	16	0	12	0	0
2	2	16	0	12	3	0
2	3	16	0	12	0	0
2	4	16	0	12	1	0
2	5	16	0	12	1	0
2	6	16	0	12	0	0
2	7	16	0	12	2	0
2	8	16	0	12	2	0
2	9	16	0	12	0	0
2	A	16	0	12	1	0
2	B	16	0	12	2	0
2	C	16	0	12	2	0
2	D	16	0	12	1	0
2	E	16	0	12	1	0
2	F	16	0	12	2	0
2	G	16	0	12	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	H	16	0	12	2	0
2	I	16	0	12	2	0
2	J	16	0	12	2	0
2	K	16	0	12	0	0
2	L	16	0	12	1	0
2	M	16	0	12	2	0
2	N	16	0	12	1	0
2	O	16	0	12	2	0
2	P	16	0	12	1	0
2	Q	16	0	12	2	0
2	R	16	0	12	1	0
2	S	16	0	12	0	0
2	T	16	0	12	2	0
2	U	16	0	12	1	0
2	V	16	0	12	0	0
2	W	16	0	12	0	0
2	X	16	0	12	2	0
2	Y	16	0	12	2	0
2	Z	16	0	12	0	0
2	a	16	0	12	0	0
2	b	16	0	12	0	0
2	c	16	0	12	0	0
2	d	16	0	12	0	0
2	e	16	0	12	0	0
2	f	16	0	12	0	0
2	g	16	0	12	0	0
2	h	16	0	12	0	0
2	i	16	0	12	0	0
2	j	16	0	12	0	0
2	k	16	0	12	0	0
2	l	16	0	12	0	0
2	m	16	0	12	0	0
3	1	9	0	0	0	0
3	2	9	0	0	0	0
3	3	9	0	0	0	0
3	4	9	0	0	0	0
3	5	9	0	0	0	0
3	6	9	0	0	0	0
3	7	9	0	0	0	0
3	8	9	0	0	0	0
3	9	9	0	0	0	0
3	A	9	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	9	0	0	0	0
3	C	9	0	0	0	0
3	D	9	0	0	0	0
3	E	9	0	0	0	0
3	F	9	0	0	0	0
3	G	9	0	0	0	0
3	H	9	0	0	0	0
3	I	9	0	0	0	0
3	J	9	0	0	0	0
3	K	9	0	0	0	0
3	L	9	0	0	0	0
3	M	9	0	0	0	0
3	N	9	0	0	0	0
3	O	9	0	0	0	0
3	P	9	0	0	0	0
3	Q	9	0	0	0	0
3	R	9	0	0	0	0
3	S	9	0	0	0	0
3	T	9	0	0	0	0
3	U	9	0	0	0	0
3	V	9	0	0	0	0
3	W	9	0	0	0	0
3	X	9	0	0	0	0
3	Y	9	0	0	0	0
3	Z	9	0	0	0	0
3	a	9	0	0	0	0
3	b	9	0	0	0	0
3	c	9	0	0	0	0
3	d	9	0	0	0	0
3	e	9	0	0	0	0
3	f	9	0	0	0	0
3	g	9	0	0	0	0
3	h	9	0	0	0	0
3	i	9	0	0	0	0
3	j	9	0	0	0	0
3	k	9	0	0	0	0
3	l	9	0	0	0	0
3	m	9	0	0	0	0
4	1	1	0	0	0	0
4	2	1	0	0	0	0
4	3	1	0	0	0	0
4	4	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	5	1	0	0	0	0
4	6	1	0	0	0	0
4	7	1	0	0	0	0
4	8	1	0	0	0	0
4	9	1	0	0	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
4	D	1	0	0	0	0
4	E	1	0	0	0	0
4	F	1	0	0	0	0
4	G	1	0	0	0	0
4	H	1	0	0	0	0
4	I	1	0	0	0	0
4	J	1	0	0	0	0
4	K	1	0	0	0	0
4	L	1	0	0	0	0
4	M	1	0	0	0	0
4	N	1	0	0	0	0
4	O	1	0	0	0	0
4	P	1	0	0	0	0
4	Q	1	0	0	0	0
4	R	1	0	0	0	0
4	S	1	0	0	0	0
4	T	1	0	0	0	0
4	U	1	0	0	0	0
4	V	1	0	0	0	0
4	W	1	0	0	0	0
4	X	1	0	0	0	0
4	Y	1	0	0	0	0
4	Z	1	0	0	0	0
4	a	1	0	0	0	0
4	b	1	0	0	0	0
4	c	1	0	0	0	0
4	d	1	0	0	0	0
4	e	1	0	0	0	0
4	f	1	0	0	0	0
4	g	1	0	0	0	0
4	h	1	0	0	0	0
4	i	1	0	0	0	0
4	j	1	0	0	0	0
4	k	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	l	1	0	0	0	0
4	m	1	0	0	0	0
5	e	5	0	0	0	0
5	l	5	0	0	0	0
6	1	120	0	0	7	0
6	2	139	0	0	1	0
6	3	102	0	0	1	0
6	4	85	0	0	3	0
6	5	109	0	0	0	0
6	6	93	0	0	2	0
6	7	69	0	0	1	0
6	8	136	0	0	3	0
6	9	101	0	0	0	0
6	A	125	0	0	2	0
6	B	96	0	0	3	0
6	C	67	0	0	0	0
6	D	133	0	0	1	0
6	E	111	0	0	7	0
6	F	86	0	0	2	0
6	G	141	0	0	4	0
6	H	123	0	0	2	0
6	I	84	0	0	0	0
6	J	116	0	0	0	0
6	K	95	0	0	1	0
6	L	70	0	0	1	0
6	M	124	0	0	3	0
6	N	98	0	0	2	0
6	O	68	0	0	2	0
6	P	142	0	0	5	0
6	Q	103	0	0	3	0
6	R	75	0	0	6	0
6	S	133	0	0	0	0
6	T	90	0	0	1	0
6	U	62	0	0	0	0
6	V	118	0	0	3	0
6	W	97	0	0	1	0
6	X	70	0	0	1	0
6	Y	94	0	0	0	0
6	Z	60	0	0	1	0
6	a	70	0	0	0	0
6	b	127	0	0	0	0
6	c	105	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	d	63	0	0	0	0
6	e	128	0	0	0	0
6	f	92	0	0	0	0
6	g	74	0	0	0	0
6	h	143	0	0	0	0
6	i	99	0	0	0	0
6	j	63	0	0	0	0
6	k	109	0	0	0	0
6	l	69	0	0	0	0
6	m	49	0	0	0	0
All	All	60911	0	55378	494	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 5.

All (494) 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:Q:119:MET:CE	1:R:115:ILE:HG12	1.64	1.25
1:M:119:MET:CE	1:O:115:ILE:HG12	1.67	1.23
1:Q:2:THR:HG1	1:V:3:MET:N	1.43	1.16
1:M:115:ILE:HG12	1:N:119:MET:CE	1.77	1.15
1:M:119:MET:HE3	1:O:115:ILE:HG12	1.23	1.15
1:E:99:LEU:HD22	1:F:97:TYR:CE2	1.81	1.13
1:8:35:THR:HG23	1:8:103:GLU:HG3	1.34	1.09
1:W:19:LYS:HE2	6:W:2019:HOH:O	1.49	1.09
1:A:105:LYS:HE2	6:A:2134:HOH:O	1.51	1.09
1:3:16:ARG:HG2	6:3:2017:HOH:O	1.52	1.08
1:A:2:THR:HG1	1:J:3:MET:N	1.52	1.06
1:A:99:LEU:HD23	1:C:97:TYR:CZ	1.90	1.06
1:2:35:THR:HG23	1:2:103:GLU:HG3	1.36	1.04
1:3:35:THR:HG23	1:3:103:GLU:HG3	1.39	1.04
1:6:35:THR:HG23	1:6:103:GLU:HG3	1.38	1.04
1:M:115:ILE:HG12	1:N:119:MET:HE3	1.35	1.03
1:1:99:LEU:HD22	1:Y:97:TYR:CE2	1.93	1.03
1:B:97:TYR:CE2	1:C:99:LEU:HD22	1.92	1.03
1:Q:119:MET:HE3	1:R:115:ILE:CG1	1.88	1.02
1:1:99:LEU:HD23	1:Y:97:TYR:CZ	1.95	1.01
1:1:119[A]:MET:CE	1:Z:115:ILE:HG12	1.90	1.01
1:A:119[A]:MET:HE3	1:B:115:ILE:HG12	1.44	1.00
1:1:99:LEU:CD2	1:Y:97:TYR:CZ	2.44	1.00

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:99:LEU:CD2	1:C:97:TYR:CE2	2.46	0.98
1:5:35:THR:CG2	1:5:103:GLU:HG3	1.94	0.97
1:7:35:THR:HG23	1:7:103:GLU:HG3	1.43	0.97
1:Q:119:MET:HE3	1:R:115:ILE:HG12	0.99	0.97
1:1:119[A]:MET:HE3	1:Z:115:ILE:CG1	1.93	0.96
1:S:35:THR:HG23	1:S:103:GLU:HG3	1.44	0.96
1:B:97:TYR:CZ	1:C:99:LEU:CD2	2.49	0.96
1:B:97:TYR:CE2	1:C:99:LEU:CD2	2.50	0.93
1:B:97:TYR:CZ	1:C:99:LEU:HD23	2.03	0.93
1:F:2:THR:HG21	1:R:121:ALA:O	1.67	0.93
1:E:143:THR:HG21	6:E:2230:HOH:O	1.67	0.93
1:A:97:TYR:CE2	1:B:99:LEU:HD22	2.05	0.92
1:1:99:LEU:CD2	1:Y:97:TYR:CE2	2.53	0.92
6:B:2162:HOH:O	1:C:143:THR:HG21	1.69	0.92
1:1:119[A]:MET:HE3	1:Z:115:ILE:HG12	0.96	0.92
1:A:99:LEU:HD23	1:C:97:TYR:CE2	2.05	0.92
1:P:143:THR:HG21	6:P:2136:HOH:O	1.70	0.92
1:D:97:TYR:CZ	1:F:99:LEU:HD22	2.04	0.91
1:I:67:TYR:CE1	1:I:99:LEU:HD23	3.73	0.91
1:D:99:LEU:HD13	1:E:97:TYR:CE2	2.07	0.89
1:E:99:LEU:CD2	1:F:97:TYR:CZ	2.54	0.89
1:D:119[A]:MET:HE3	1:F:115:ILE:HG12	1.56	0.88
1:J:143:THR:HG23	1:J:144:LYS:HD3	2.00	0.88
1:Z:67:TYR:CE1	1:Z:99:LEU:HD23	2.08	0.87
1:A:97:TYR:CZ	1:B:99:LEU:CD2	2.57	0.87
1:M:115:ILE:HG12	1:N:119:MET:HE1	1.54	0.87
1:9:99:LEU:HD13	1:A:97:TYR:CE2	112.73	0.86
1:2:119[A]:MET:HE3	1:4:115:ILE:HG12	1.56	0.85
1:5:99:LEU:HD13	1:7:97:TYR:CE2	2.10	0.85
1:1:35:THR:HG22	1:1:103:GLU:HG3	1.57	0.85
1:8:67:TYR:CE1	1:8:99:LEU:HD12	2.12	0.85
1:A:99:LEU:CD2	1:C:97:TYR:CZ	2.61	0.83
1:I:67:TYR:HE1	1:I:99:LEU:HD23	3.06	0.83
1:2:35:THR:HG23	1:2:103:GLU:CG	2.10	0.82
1:5:35:THR:HG22	1:5:103:GLU:HG3	1.62	0.82
1:R:6:LYS:HE2	6:R:2003:HOH:O	1.79	0.82
1:H:67:TYR:CE1	1:H:99:LEU:HD23	2.14	0.82
1:A:97:TYR:CZ	1:B:99:LEU:HD22	2.15	0.81
1:S:35:THR:CG2	1:S:103:GLU:HG3	2.09	0.81
1:7:35:THR:CG2	1:7:103:GLU:HG3	2.11	0.81
1:F:119[A]:MET:HE3	1:G:115:ILE:HG12	47.66	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:V:2080:HOH:O	1:W:129:HIS:HB3	1.82	0.80
1:M:119:MET:HE1	1:O:115:ILE:HG12	1.62	0.80
6:1:2014:HOH:O	1:G:16:ARG:HG2	132.61	0.80
1:8:67:TYR:HE1	1:8:99:LEU:HD12	1.47	0.79
1:Z:67:TYR:HE1	1:Z:99:LEU:HD23	1.44	0.79
1:5:99:LEU:HD13	1:7:97:TYR:CZ	2.18	0.78
1:1:35:THR:HG21	6:1:2086:HOH:O	1.84	0.78
1:7:35:THR:HG23	1:7:103:GLU:CG	2.12	0.78
1:6:35:THR:HG23	1:6:103:GLU:CG	2.14	0.78
1:G:119:MET:CE	1:H:115:ILE:HG12	2.13	0.78
1:E:99:LEU:CD2	1:F:97:TYR:CE2	2.64	0.78
1:M:119:MET:HE3	1:O:115:ILE:CG1	2.11	0.78
1:E:99:LEU:HD22	1:F:97:TYR:CZ	2.15	0.78
1:2:138:LEU:HD22	1:4:43:LEU:HD11	1.64	0.77
1:I:119[A]:MET:HE3	1:J:115:ILE:HG12	51.39	0.77
1:B:143:THR:HG23	1:B:144:LYS:HD3	2.02	0.77
1:S:143:THR:HG22	1:S:144:LYS:HD3	1.66	0.77
1:5:35:THR:HG23	1:5:103:GLU:HG3	1.67	0.77
1:Z:67:TYR:CE1	1:Z:99:LEU:CD2	2.67	0.77
1:Z:67:TYR:HE1	1:Z:99:LEU:CD2	1.98	0.77
1:S:35:THR:HG23	1:S:103:GLU:CG	2.13	0.76
1:P:119[A]:MET:HE3	1:Q:115:ILE:HG12	1.68	0.76
1:D:67:TYR:CE1	1:D:99:LEU:HD12	2.22	0.75
1:G:143:THR:HG23	1:G:144[A]:LYS:HD3	1.96	0.75
1:M:43:LEU:HD11	1:N:138:LEU:HD22	1.66	0.75
1:E:143:THR:HG23	1:E:144:LYS:HD3	2.20	0.75
1:H:12:GLU:H	1:H:12:GLU:CD	1.89	0.75
1:O:12:GLU:CD	1:O:12:GLU:H	1.88	0.75
1:S:115:ILE:HG12	1:U:119[A]:MET:HE3	1.67	0.74
1:X:143:THR:HG22	6:X:2070:HOH:O	1.86	0.74
1:Z:37:LYS:HB2	1:Z:40:GLU:CD	2.08	0.74
1:8:129:HIS:ND1	6:8:2117:HOH:O	2.22	0.73
1:W:43:LEU:HD11	1:X:138:LEU:HD22	1.71	0.73
1:P:43:LEU:HD11	1:R:138:LEU:HD22	1.69	0.73
1:1:35:THR:CG2	1:1:103:GLU:HG3	2.19	0.72
1:I:119[A]:MET:CE	1:J:115:ILE:HG12	51.17	0.72
1:R:12:GLU:H	1:R:12:GLU:CD	1.93	0.72
1:6:35:THR:CG2	1:6:103:GLU:HG3	2.19	0.72
1:5:35:THR:HG23	1:5:103:GLU:CG	2.20	0.71
1:9:99:LEU:HD13	1:A:97:TYR:CZ	113.90	0.71
1:5:97:TYR:CZ	1:6:99:LEU:HD23	2.26	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:99:LEU:HD22	1:C:97:TYR:CE2	2.23	0.71
1:D:67:TYR:HE1	1:D:99:LEU:HD12	1.54	0.71
1:E:138:LEU:HD22	1:F:43:LEU:HD11	28.82	0.71
1:R:143:THR:HG21	6:R:2073:HOH:O	1.91	0.71
6:8:2098:HOH:O	1:9:129:HIS:ND1	2.24	0.71
1:4:106:LYS:NZ	6:4:2006:HOH:O	2.24	0.70
1:A:143:THR:HG23	1:A:144:LYS:HD3	1.96	0.70
1:5:99:LEU:CD1	1:7:97:TYR:CE2	2.75	0.69
1:D:43:LEU:HD11	1:E:138:LEU:HD22	1.74	0.69
1:G:99:LEU:HD23	1:I:97:TYR:CE2	2.27	0.69
1:B:12:GLU:OE2	6:B:2014:HOH:O	2.09	0.69
1:D:143:THR:HG23	1:D:144:LYS:HD3	2.08	0.69
1:2:99:LEU:HD22	1:3:97:TYR:CZ	2.28	0.68
1:M:143:THR:HG23	1:M:144:LYS:HD3	2.06	0.68
1:Q:106:LYS:HE2	6:Q:2077:HOH:O	1.93	0.68
1:U:12:GLU:CD	1:U:12:GLU:H	1.97	0.68
1:F:2:THR:CG2	1:P:3:MET:N	2.57	0.68
1:B:138:LEU:HD22	1:C:43:LEU:HD11	1.74	0.68
1:F:119[A]:MET:CE	1:G:115:ILE:HG12	48.35	0.68
1:H:67:TYR:HE1	1:H:99:LEU:HD23	1.56	0.67
1:A:115:ILE:HG12	1:C:119:MET:CE	2.24	0.67
1:G:119:MET:HE3	1:H:115:ILE:HG12	1.74	0.67
1:8:99:LEU:HD13	1:9:97:TYR:CE2	2.30	0.67
1:1:3:MET:HA	6:1:2001:HOH:O	1.94	0.67
1:A:119[A]:MET:CE	1:B:115:ILE:HG12	2.23	0.67
1:A:138:LEU:HD22	1:B:43:LEU:HD11	1.75	0.67
1:I:67:TYR:HE1	1:I:99:LEU:CD2	3.26	0.67
1:J:43:LEU:HD11	1:K:138:LEU:HD22	1.77	0.67
1:S:143:THR:CG2	1:S:144:LYS:HD3	2.24	0.66
1:E:16:ARG:HG2	6:E:2025:HOH:O	1.96	0.66
1:G:3:MET:HE3	6:G:2045:HOH:O	1.95	0.66
1:N:115:ILE:HG12	1:O:119:MET:CE	2.24	0.66
1:A:115:ILE:HG12	1:C:119:MET:HE1	1.77	0.66
1:M:115:ILE:CG1	1:N:119:MET:HE3	2.19	0.66
1:G:143:THR:HG22	6:G:2204:HOH:O	1.95	0.66
1:I:38:LYS:HG3	1:I:99:LEU:O	2.75	0.66
1:S:138:LEU:HD22	1:T:43:LEU:HD11	1.78	0.66
1:F:143:THR:HG23	1:F:144:LYS:HD3	2.27	0.66
1:S:115:ILE:HG12	1:U:119[A]:MET:CE	2.24	0.66
1:K:138:LEU:HD22	1:L:43:LEU:HD11	28.63	0.66
1:M:138:LEU:HD22	1:O:43:LEU:HD11	1.77	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:V:43:LEU:HD11	1:W:138:LEU:HD22	1.78	0.65
1:5:35:THR:CG2	1:5:103:GLU:CG	2.72	0.65
6:1:2077:HOH:O	1:Y:129:HIS:CG	2.48	0.65
1:9:143:THR:HG23	1:9:144[A]:LYS:HD3	1.79	0.65
1:L:143:THR:HG23	1:L:144:LYS:HD3	2.21	0.65
1:I:67:TYR:CE1	1:I:99:LEU:CD2	3.98	0.64
1:Y:43:LEU:HD11	1:Z:138:LEU:HD22	1.80	0.64
1:7:67:TYR:CE1	1:7:99:LEU:HD12	2.32	0.64
1:3:43:LEU:HD11	1:4:138:LEU:HD22	1.80	0.64
1:5:43:LEU:HD11	1:7:138:LEU:HD22	1.78	0.64
1:D:119[A]:MET:CE	1:F:115:ILE:HG12	2.27	0.64
1:T:117:LYS:HE2	6:T:2070:HOH:O	1.98	0.64
1:L:138:LEU:HD22	1:M:43:LEU:HD11	98.99	0.64
1:N:43:LEU:HD11	1:O:138:LEU:HD22	1.80	0.64
1:D:138:LEU:HD22	1:F:43:LEU:HD11	1.80	0.63
1:8:138:LEU:HD22	1:A:43:LEU:HD11	104.30	0.63
1:S:43:LEU:HD11	1:U:138:LEU:HD22	1.80	0.63
1:D:12:GLU:H	1:D:12:GLU:CD	2.18	0.63
1:K:37:LYS:O	1:K:40:GLU:HG3	1.98	0.63
1:Q:2:THR:OG1	1:V:3:MET:N	2.24	0.63
1:6:97:TYR:CE2	1:7:99:LEU:HD13	2.34	0.62
1:Z:38:LYS:HG3	1:Z:99:LEU:O	1.99	0.62
1:6:138:LEU:HD22	1:7:43:LEU:HD11	1.81	0.62
1:P:138:LEU:HD22	1:Q:43:LEU:HD11	1.81	0.62
1:F:2:THR:CG2	1:R:121:ALA:O	2.44	0.62
1:G:43:LEU:HD11	1:I:138:LEU:HD22	1.80	0.62
1:1:43:LEU:HD11	1:Y:138:LEU:HD22	1.81	0.62
1:F:12:GLU:CD	1:F:12:GLU:H	2.03	0.61
1:H:143:THR:HG23	1:H:144:LYS:HD3	2.26	0.61
6:M:2060:HOH:O	1:O:143:THR:HG21	1.98	0.61
1:P:143:THR:HG23	1:P:144[A]:LYS:HD3	1.82	0.61
1:2:138:LEU:CD2	1:4:43:LEU:HD11	2.30	0.61
1:2:119[A]:MET:CE	1:4:115:ILE:HG12	2.29	0.61
1:7:67:TYR:HE1	1:7:99:LEU:HD12	1.65	0.61
1:E:19:LYS:HE2	6:E:2046:HOH:O	2.00	0.61
1:5:138:LEU:HD22	1:6:43:LEU:HD11	1.82	0.61
1:G:119:MET:HE1	1:H:115:ILE:HG12	1.82	0.60
1:K:43:LEU:HD11	1:L:138:LEU:HD22	1.81	0.60
1:X:143:THR:HG23	1:X:144:LYS:HD3	1.83	0.60
1:R:143:THR:HG23	1:R:144:LYS:HD3	1.83	0.60
1:C:138:LEU:HD22	1:D:43:LEU:HD11	69.03	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:43:LEU:HD11	1:C:138:LEU:HD22	1.84	0.60
1:M:43:LEU:HD11	1:N:138:LEU:CD2	2.32	0.60
6:A:2030:HOH:O	1:J:16:ARG:HG2	2.01	0.59
1:A:2:THR:OG1	1:J:3:MET:N	2.30	0.59
1:D:115:ILE:HG12	1:E:119:MET:CE	2.33	0.59
1:8:103:GLU:HG2	6:8:2045:HOH:O	2.03	0.58
1:G:3:MET:CE	6:G:2045:HOH:O	2.51	0.58
1:H:38:LYS:HG3	1:H:99:LEU:O	2.03	0.58
1:P:43:LEU:HD11	1:R:138:LEU:CD2	2.34	0.58
1:4:117:LYS:NZ	6:4:2068:HOH:O	2.25	0.58
1:Q:138:LEU:HD22	1:R:43:LEU:HD11	1.84	0.58
1:B:97:TYR:CD2	1:C:99:LEU:HD22	2.37	0.58
1:E:43:LEU:HD11	1:G:138:LEU:HD22	47.01	0.58
1:1:138:LEU:HD22	1:Z:43:LEU:HD11	1.86	0.58
1:6:129:HIS:HB3	6:6:2087:HOH:O	2.03	0.57
1:T:138:LEU:HD22	1:U:43:LEU:HD11	1.85	0.57
1:D:38:LYS:HG3	1:D:99:LEU:O	2.05	0.57
1:2:127:VAL:HG12	1:A:21:GLU:HB2	112.28	0.57
1:E:67:TYR:CE1	1:E:99:LEU:HD13	2.40	0.57
1:I:138:LEU:HD22	1:J:43:LEU:HD11	87.29	0.57
1:Q:119:MET:HE1	1:R:115:ILE:HG12	1.78	0.57
1:D:115:ILE:HG12	1:E:119:MET:HE1	1.87	0.57
1:J:115:ILE:HG12	1:K:119:MET:CE	2.35	0.56
1:H:138:LEU:HD22	1:I:43:LEU:HD11	1.86	0.56
1:N:43:LEU:HD11	1:O:138:LEU:CD2	2.35	0.56
1:3:35:THR:HG23	1:3:103:GLU:CG	2.26	0.56
1:R:117:LYS:HD3	6:R:2018:HOH:O	2.04	0.56
1:8:3:MET:HE2	1:8:53:PRO:CB	2.35	0.56
1:1:99:LEU:HD22	1:Y:97:TYR:CD2	2.37	0.56
1:F:85:TYR:CD2	2:F:1145:DUR:H2'2	2.41	0.56
1:A:97:TYR:CZ	1:B:99:LEU:HD23	2.41	0.56
1:Y:115:ILE:HG12	1:Z:119[A]:MET:HE3	1.86	0.56
1:Z:143:THR:HG23	1:Z:144[B]:LYS:HD3	1.86	0.56
1:8:143:THR:HG23	1:8:144:LYS:HD3	1.86	0.56
1:M:128:GLU:OE1	1:M:129:HIS:CD2	2.58	0.56
1:T:37:LYS:HB2	1:T:40:GLU:CD	2.26	0.56
1:B:37:LYS:HB2	1:B:40:GLU:CD	3.01	0.56
1:K:43:LEU:HD11	1:M:138:LEU:HD22	84.02	0.56
1:9:119[A]:MET:HE1	1:9:122:VAL:HG11	1.88	0.55
1:F:138:LEU:HD22	1:G:43:LEU:HD11	56.20	0.55
1:2:43:LEU:HD11	1:3:138:LEU:HD22	1.87	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:9:67:TYR:CE1	1:9:99:LEU:HD12	2.41	0.55
1:W:119:MET:CE	1:W:122:VAL:HG11	2.37	0.55
1:A:12:GLU:CD	1:A:12:GLU:H	2.21	0.54
1:J:11:ASP:HB2	1:J:12:GLU:OE2	4.08	0.54
1:8:35:THR:HG23	1:8:103:GLU:CG	2.24	0.54
1:1:99:LEU:HD23	1:Y:97:TYR:OH	2.06	0.54
1:A:99:LEU:CD2	1:C:97:TYR:CD2	2.91	0.54
1:H:143:THR:CG2	1:H:144:LYS:HD3	3.09	0.54
1:6:119[A]:MET:CE	1:7:115:ILE:HG12	2.37	0.54
1:B:115:ILE:HG12	1:D:119[A]:MET:CE	34.60	0.54
1:Z:106:LYS:NZ	6:Z:2004:HOH:O	2.40	0.54
1:9:119[A]:MET:CE	1:9:122:VAL:CG1	2.86	0.54
6:E:2047:HOH:O	1:F:119[A]:MET:HG2	2.08	0.54
1:7:100:ARG:NE	6:7:2043:HOH:O	2.24	0.53
1:F:2:THR:HG21	1:P:3:MET:N	2.22	0.53
1:D:85:TYR:CD2	2:D:1145:DUR:H2'2	2.60	0.53
1:W:43:LEU:HD11	1:X:138:LEU:CD2	2.38	0.53
1:Z:143:THR:HG23	1:Z:144[A]:LYS:HD3	1.90	0.53
1:H:97:TYR:CE2	1:I:99:LEU:HD23	2.43	0.53
1:H:143:THR:HG22	6:H:2257:HOH:O	2.07	0.53
1:N:115:ILE:HG12	1:O:119:MET:HE1	1.91	0.53
1:P:119[A]:MET:CE	1:Q:115:ILE:HG12	2.38	0.53
1:H:143:THR:HG23	1:H:144:LYS:CD	2.93	0.53
1:Q:143:THR:HG23	1:Q:144:LYS:HD3	1.91	0.53
1:5:67:TYR:CE1	1:5:99:LEU:HD12	2.44	0.53
1:E:100:ARG:NE	6:E:2121:HOH:O	2.39	0.53
1:K:12:GLU:HG3	6:K:2017:HOH:O	2.08	0.53
1:V:138:LEU:HD22	1:X:43:LEU:HD11	1.90	0.53
1:K:7:ILE:HD11	1:K:48:VAL:CG1	2.86	0.52
1:1:35:THR:CG2	1:1:103:GLU:CG	2.86	0.52
1:9:43:LEU:HD11	1:A:138:LEU:HD22	126.66	0.52
1:N:115:ILE:HG12	1:O:119:MET:HE3	1.91	0.52
1:K:6:LYS:HZ3	1:K:51:GLU:CD	4.34	0.52
1:C:85:TYR:CG	2:C:1145:DUR:H2'2	2.44	0.52
6:1:2077:HOH:O	1:Y:129:HIS:HB3	2.08	0.52
1:Z:37:LYS:HB2	1:Z:40:GLU:OE2	2.07	0.52
1:C:85:TYR:CD2	2:C:1145:DUR:H2'2	2.45	0.52
1:F:35:THR:O	1:F:42:LYS:NZ	3.29	0.52
1:G:58:ALA:HB3	1:G:81:ILE:HB	2.04	0.52
1:H:119:MET:HE3	1:H:122:VAL:HG13	1.92	0.52
1:E:43:LEU:HD11	1:F:138:LEU:HD22	1.92	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:9:67:TYR:HE1	1:9:99:LEU:HD12	1.76	0.51
1:P:35:THR:CG2	6:P:2049:HOH:O	2.58	0.51
1:K:11:ASP:OD1	1:K:13:THR:N	2.64	0.51
1:L:4:GLN:NE2	6:L:2005:HOH:O	2.44	0.51
1:K:76:ASN:O	1:K:77:SER:HB2	2.42	0.51
1:S:97:TYR:CE2	1:T:99:LEU:HD23	2.46	0.51
1:A:99:LEU:HD22	1:C:97:TYR:CD2	2.46	0.51
1:A:138:LEU:CD2	1:B:43:LEU:HD11	2.41	0.51
1:K:6:LYS:NZ	1:K:51:GLU:CD	5.01	0.50
1:Z:67:TYR:CE1	1:Z:99:LEU:HD22	2.44	0.50
1:K:138:LEU:CD2	1:L:43:LEU:HD11	28.22	0.50
1:K:99:LEU:HD23	1:L:97:TYR:CE2	2.46	0.50
1:Y:115:ILE:HG12	1:Z:119[A]:MET:CE	2.41	0.50
1:9:119[A]:MET:CE	1:9:122:VAL:HG11	2.42	0.50
1:X:85:TYR:CD2	2:X:1145:DUR:H2'2	2.47	0.50
1:B:115:ILE:HG12	1:D:119[A]:MET:HE3	33.76	0.49
1:J:138:LEU:HD22	1:L:43:LEU:HD11	1.94	0.49
1:G:138:LEU:HD22	1:H:43:LEU:HD11	1.93	0.49
1:P:105:LYS:HE3	6:P:2046:HOH:O	2.12	0.49
1:Q:97:TYR:CE2	1:R:99:LEU:HD23	2.47	0.49
1:9:43:LEU:HD11	1:A:138:LEU:CD2	127.46	0.49
1:K:7:ILE:HD11	1:K:48:VAL:HG11	2.46	0.49
1:E:99:LEU:HD23	1:F:97:TYR:CZ	2.46	0.49
1:B:138:LEU:CD2	1:C:43:LEU:HD11	2.43	0.49
1:F:38:LYS:HG3	1:F:99:LEU:O	2.12	0.49
1:J:43:LEU:HD11	1:K:138:LEU:CD2	2.41	0.49
1:O:100:ARG:NH1	1:O:101:ASP:O	2.46	0.49
2:2:1145:DUR:O5'	2:2:1145:DUR:H6	2.12	0.49
1:B:85:TYR:CD2	2:B:1145:DUR:H2'2	2.47	0.49
1:G:99:LEU:HD23	1:I:97:TYR:CD2	2.48	0.49
1:P:97:TYR:CE2	1:Q:99:LEU:HD23	2.47	0.49
1:D:138:LEU:CD2	1:F:43:LEU:HD11	2.43	0.48
1:E:138:LEU:CD2	1:F:43:LEU:HD11	28.36	0.48
1:Q:101:ASP:HB3	6:Q:2074:HOH:O	2.12	0.48
1:S:138:LEU:CD2	1:T:43:LEU:HD11	2.43	0.48
1:H:85:TYR:CG	2:H:1145:DUR:H2'2	2.48	0.48
1:M:85:TYR:CD2	2:M:1145:DUR:H2'2	2.51	0.48
1:J:143:THR:CG2	1:J:144:LYS:HD3	2.97	0.48
1:T:97:TYR:CE2	1:U:99:LEU:HD23	2.49	0.48
1:B:38:LYS:HG3	1:B:99:LEU:O	2.14	0.48
1:I:127:VAL:HG12	1:X:21:GLU:HB2	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:12:GLU:H	1:K:12:GLU:CD	2.14	0.48
1:O:85:TYR:CG	2:O:1145:DUR:H2'2	2.49	0.48
1:8:43:LEU:HD11	1:9:138:LEU:HD22	1.94	0.48
1:B:43:LEU:HD11	1:D:138:LEU:HD22	44.55	0.48
1:J:85:TYR:CD2	2:J:1145:DUR:H2'2	2.66	0.48
1:5:43:LEU:HD11	1:7:138:LEU:CD2	2.44	0.48
1:V:58:ALA:HB3	1:V:81:ILE:HB	1.95	0.48
1:G:35:THR:HG22	1:G:103:GLU:HG3	1.96	0.47
1:8:3:MET:HE2	1:8:53:PRO:HB3	1.95	0.47
1:8:67:TYR:HE1	1:8:99:LEU:CD1	2.22	0.47
1:1:43:LEU:HD11	1:Y:138:LEU:CD2	2.45	0.47
6:F:2028:HOH:O	1:P:16:ARG:HG2	2.14	0.47
1:3:99:LEU:HD22	1:4:97:TYR:CZ	2.50	0.47
1:8:119[A]:MET:CE	1:A:115:ILE:HG12	107.46	0.47
1:8:138:LEU:CD2	1:A:43:LEU:HD11	104.62	0.47
1:B:85:TYR:CG	2:B:1145:DUR:H2'2	2.49	0.47
1:M:99:LEU:HD23	1:N:97:TYR:CE2	2.50	0.47
1:Q:117:LYS:HE2	6:Q:2024:HOH:O	2.14	0.47
1:K:38:LYS:HE2	1:K:99:LEU:O	2.45	0.47
1:B:97:TYR:OH	1:C:99:LEU:HD23	2.13	0.47
1:P:138:LEU:CD2	1:Q:43:LEU:HD11	2.45	0.47
1:D:99:LEU:HD13	1:E:97:TYR:CD2	2.47	0.47
1:K:98:ALA:HB1	1:K:100:ARG:O	2.38	0.47
1:5:99:LEU:CD1	1:7:97:TYR:CD2	2.98	0.47
1:K:105:LYS:HA	1:K:105:LYS:HD2	1.79	0.47
1:M:12:GLU:H	1:M:12:GLU:CD	2.18	0.47
1:7:85:TYR:CG	2:7:1145:DUR:H2'2	2.50	0.46
1:D:43:LEU:HD11	1:E:138:LEU:CD2	2.44	0.46
1:F:85:TYR:CG	2:F:1145:DUR:H2'2	2.50	0.46
1:Q:37:LYS:HB2	1:Q:40:GLU:OE1	2.14	0.46
1:5:38:LYS:HG3	1:5:99:LEU:O	2.15	0.46
1:8:58:ALA:HB3	1:8:81:ILE:HB	1.96	0.46
1:1:117:LYS:HE3	6:1:2054:HOH:O	2.16	0.46
1:5:35:THR:HG23	1:5:103:GLU:HG2	1.94	0.46
1:M:35:THR:HG23	6:M:2053:HOH:O	2.15	0.46
1:6:119[A]:MET:HE3	1:7:115:ILE:HG12	1.95	0.46
2:8:1145:DUR:O5'	2:8:1145:DUR:H6	2.15	0.46
1:J:143:THR:HG23	1:J:144:LYS:CD	2.73	0.46
1:K:27:ASP:OD2	1:K:109:ARG:HD2	2.47	0.46
1:Y:85:TYR:CG	2:Y:1145:DUR:H2'2	2.50	0.46
1:6:119[A]:MET:HE3	1:6:119[A]:MET:HB2	1.75	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:7:38:LYS:HG3	1:7:99:LEU:O	2.16	0.46
1:E:43:LEU:HD11	1:G:138:LEU:CD2	47.82	0.46
1:N:103:GLU:HG2	6:N:2028:HOH:O	2.15	0.46
6:H:2216:HOH:O	1:K:19:LYS:HE3	2.14	0.46
1:Q:85:TYR:CD2	2:Q:1145:DUR:H2'2	2.51	0.46
1:Y:43:LEU:HD11	1:Z:138:LEU:CD2	2.44	0.46
1:8:99:LEU:HD13	1:9:97:TYR:CD2	2.50	0.46
1:9:38:LYS:HG3	1:9:99:LEU:O	2.15	0.46
1:D:99:LEU:HD13	1:E:97:TYR:CZ	2.47	0.46
1:8:85:TYR:CG	2:8:1145:DUR:H2'2	2.51	0.46
1:O:85:TYR:CD2	2:O:1145:DUR:H2'2	2.51	0.46
1:9:143:THR:HG23	1:9:144[B]:LYS:HD3	1.98	0.46
1:L:138:LEU:CD2	1:M:43:LEU:HD11	99.03	0.46
1:V:43:LEU:HD11	1:W:138:LEU:CD2	2.45	0.46
1:V:97:TYR:CE2	1:X:99:LEU:HD23	2.51	0.46
1:U:12:GLU:CD	1:U:12:GLU:N	2.68	0.46
1:5:127:VAL:HG12	1:Z:21:GLU:HB2	1.98	0.45
1:N:99:LEU:HD23	1:O:97:TYR:CE2	2.51	0.45
1:L:21:GLU:HB2	1:T:127:VAL:HG12	1.98	0.45
1:M:100:ARG:NE	6:M:2100:HOH:O	2.21	0.45
1:O:58:ALA:HB3	1:O:81:ILE:HB	1.98	0.45
1:V:3:MET:HA	6:V:2002:HOH:O	2.15	0.45
1:6:138:LEU:CD2	1:7:43:LEU:HD11	2.45	0.45
1:X:85:TYR:CG	2:X:1145:DUR:H2'2	2.52	0.45
1:B:35:THR:HG22	1:B:103:GLU:CG	2.76	0.45
1:C:37:LYS:O	1:C:40:GLU:HB2	2.51	0.45
1:E:119:MET:HE3	1:E:119:MET:HB2	1.37	0.45
1:I:85:TYR:CG	2:I:1145:DUR:H2'2	2.50	0.45
1:J:115:ILE:HG12	1:K:119:MET:HE3	1.98	0.45
1:1:103:GLU:HG2	6:1:2087:HOH:O	2.15	0.45
1:E:85:TYR:CG	2:E:1145:DUR:H2'2	2.68	0.45
1:K:17:ILE:HG21	1:K:28:LEU:HD23	2.42	0.45
1:E:37:LYS:HB2	1:E:40:GLU:CD	2.73	0.45
1:W:119:MET:CE	1:W:122:VAL:CG1	2.95	0.45
1:E:127:VAL:HG12	1:H:21:GLU:HB2	1.99	0.44
1:9:99:LEU:CD1	1:A:97:TYR:CE2	113.69	0.44
1:F:10:LEU:O	6:F:2011:HOH:O	2.20	0.44
1:I:85:TYR:CD2	2:I:1145:DUR:H2'2	2.58	0.44
1:I:143:THR:HG23	1:I:144:LYS:HD3	2.07	0.44
1:J:3:MET:HE2	1:J:53:PRO:HB3	2.00	0.44
1:1:99:LEU:CD2	1:Y:97:TYR:CE1	2.95	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:100:ARG:NH2	6:E:2121:HOH:O	2.49	0.44
1:F:119[A]:MET:HE3	1:G:115:ILE:H	48.59	0.44
1:K:43:LEU:HD11	1:L:138:LEU:CD2	2.48	0.44
1:2:35:THR:CG2	1:2:103:GLU:OE2	2.65	0.44
1:4:3:MET:HA	6:4:2001:HOH:O	2.18	0.44
1:G:85:TYR:CD2	2:G:1145:DUR:H2'2	2.52	0.44
6:G:2147:HOH:O	1:I:129:HIS:CG	2.71	0.44
1:J:3:MET:HB3	1:J:3:MET:HE2	1.76	0.44
1:I:138:LEU:CD2	1:J:43:LEU:HD11	87.54	0.44
1:J:97:TYR:CE2	1:L:99:LEU:HD23	2.53	0.44
1:4:35:THR:HG22	1:4:103:GLU:HG2	2.00	0.44
1:8:38:LYS:HG3	1:8:99:LEU:O	2.18	0.44
1:G:43:LEU:HD11	1:I:138:LEU:CD2	2.47	0.44
1:I:119[A]:MET:HE2	1:J:115:ILE:HG12	50.66	0.44
1:T:85:TYR:CG	2:T:1145:DUR:H2'2	2.53	0.44
1:B:58:ALA:HB3	1:B:81:ILE:HB	2.02	0.44
1:8:119[A]:MET:HE3	1:A:115:ILE:HG12	107.37	0.43
1:J:85:TYR:CG	2:J:1145:DUR:H2'2	2.87	0.43
1:O:143:THR:HG22	6:O:2068:HOH:O	2.17	0.43
1:P:35:THR:HG22	6:P:2049:HOH:O	2.17	0.43
1:M:85:TYR:CG	2:M:1145:DUR:H2'2	2.60	0.43
1:R:32:GLU:OE2	6:R:2022:HOH:O	2.21	0.43
1:5:85:TYR:CD2	2:5:1145:DUR:H2'2	2.53	0.43
1:A:38:LYS:HG3	1:A:99:LEU:O	2.19	0.43
1:5:67:TYR:HE1	1:5:99:LEU:HD12	1.83	0.43
1:L:85:TYR:CD2	2:L:1145:DUR:H2'2	2.55	0.43
1:A:115:ILE:HG12	1:C:119:MET:HE3	1.98	0.43
1:H:85:TYR:CD2	2:H:1145:DUR:H2'2	2.54	0.43
1:L:119:MET:CE	1:L:122:VAL:HG11	2.48	0.43
1:U:85:TYR:CD2	2:U:1145:DUR:H2'2	2.53	0.43
1:Y:99:LEU:HD23	1:Z:97:TYR:CE2	2.54	0.43
1:V:37:LYS:O	1:V:40:GLU:HB2	2.18	0.43
1:G:85:TYR:CG	2:G:1145:DUR:H2'2	2.53	0.43
1:B:35:THR:HG22	1:B:103:GLU:HG2	2.18	0.43
1:H:119:MET:CE	1:H:122:VAL:CG1	2.97	0.43
1:K:143:THR:HG23	1:K:144:LYS:HD3	2.01	0.43
1:S:99:LEU:HD23	1:U:97:TYR:CE2	2.53	0.43
1:R:85:TYR:CD2	2:R:1145:DUR:H2'2	2.54	0.43
1:M:58:ALA:HB3	1:M:81:ILE:HB	2.02	0.42
1:N:86:LYS:NZ	6:N:2057:HOH:O	2.52	0.42
1:4:85:TYR:CD2	2:4:1145:DUR:H2'2	2.55	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:1145:DUR:H6	2:A:1145:DUR:O5'	2.18	0.42
1:Y:58:ALA:HB3	1:Y:81:ILE:HB	2.01	0.42
1:7:85:TYR:CD2	2:7:1145:DUR:H2'2	2.54	0.42
1:F:138:LEU:CD2	1:G:43:LEU:HD11	55.64	0.42
1:M:128:GLU:OE1	1:M:129:HIS:HD2	2.02	0.42
1:9:119[A]:MET:CE	1:9:122:VAL:HG13	2.49	0.42
1:D:40:GLU:OE2	6:D:2069:HOH:O	2.21	0.42
1:2:119[A]:MET:HG2	6:2:2065:HOH:O	2.19	0.42
1:2:35:THR:HG23	1:2:103:GLU:CD	2.40	0.42
1:F:67:TYR:CE1	1:F:99:LEU:HD13	2.55	0.42
1:F:2:THR:HG22	1:P:3:MET:N	2.31	0.42
6:B:2162:HOH:O	1:C:143:THR:CG2	2.44	0.42
1:4:12:GLU:CD	1:4:12:GLU:H	2.23	0.42
1:H:43:LEU:HD11	1:J:138:LEU:HD22	56.22	0.42
1:P:85:TYR:CD2	2:P:1145:DUR:H2'2	2.55	0.42
1:5:138:LEU:CD2	1:6:43:LEU:HD11	2.48	0.42
1:Q:58:ALA:HB3	1:Q:81:ILE:HB	2.01	0.42
1:R:117:LYS:CD	6:R:2018:HOH:O	2.66	0.42
1:T:37:LYS:HE2	1:T:37:LYS:HB3	1.89	0.42
1:Y:85:TYR:CD2	2:Y:1145:DUR:H2'2	2.54	0.42
1:K:58:ALA:HB3	1:K:81:ILE:HB	2.19	0.41
1:S:12:GLU:H	1:S:12:GLU:CD	2.23	0.41
1:A:97:TYR:CE1	1:B:99:LEU:CD2	3.01	0.41
1:V:99:LEU:HD23	1:W:97:TYR:CE2	2.55	0.41
1:V:117:LYS:HD2	6:V:2036:HOH:O	2.20	0.41
1:O:101:ASP:HB3	6:O:2028:HOH:O	2.20	0.41
1:R:6:LYS:CE	6:R:2003:HOH:O	2.53	0.41
1:2:85:TYR:CD2	2:2:1145:DUR:H2'2	2.56	0.41
1:J:99:LEU:HD23	1:K:97:TYR:CE2	2.55	0.41
1:P:103:GLU:HG2	6:P:2048:HOH:O	2.20	0.41
1:T:85:TYR:CD2	2:T:1145:DUR:H2'2	2.56	0.41
1:C:21:GLU:HB2	1:O:127:VAL:HG12	2.03	0.41
1:H:138:LEU:CD2	1:I:43:LEU:HD11	2.61	0.41
1:U:143:THR:HG23	1:U:144:LYS:HD3	2.02	0.41
1:B:97:TYR:CE1	1:C:99:LEU:CD2	3.01	0.41
1:H:119:MET:HE1	1:H:122:VAL:HG11	2.02	0.41
1:1:127:VAL:HG12	1:6:21:GLU:HB2	2.03	0.41
1:S:115:ILE:H	1:U:119[A]:MET:HE3	1.85	0.41
1:2:85:TYR:CG	2:2:1145:DUR:H2'2	2.56	0.41
1:D:58:ALA:HB3	1:D:81:ILE:HB	2.03	0.41
1:P:119[A]:MET:HE3	1:Q:115:ILE:H	1.86	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:Q:85:TYR:CG	2:Q:1145:DUR:H2'2	2.55	0.41
1:5:129:HIS:HB3	6:6:2062:HOH:O	2.21	0.41
1:I:122:VAL:O	1:I:122:VAL:HG23	2.42	0.41
1:M:35:THR:HG22	1:M:103:GLU:CG	2.78	0.41
1:2:35:THR:CG2	1:2:103:GLU:CD	2.90	0.40
1:8:67:TYR:CE1	1:8:99:LEU:CD1	2.94	0.40
1:D:119[A]:MET:HE3	1:F:115:ILE:H	1.85	0.40
1:M:138:LEU:CD2	1:O:43:LEU:HD11	2.47	0.40
1:1:38:LYS:HG3	1:1:99:LEU:O	2.20	0.40
1:B:21:GLU:HB2	1:F:127:VAL:HG12	2.03	0.40
1:H:12:GLU:CD	1:H:12:GLU:N	2.67	0.40
1:H:99:LEU:HA	1:H:99:LEU:HD12	1.86	0.40
1:N:85:TYR:CD2	2:N:1145:DUR:H2'2	2.56	0.40
1:8:3:MET:HE2	1:8:53:PRO:CG	2.51	0.40
1:8:21:GLU:HB2	1:D:127:VAL:HG12	113.14	0.40
1:D:97:TYR:CE2	1:F:99:LEU:HD22	2.52	0.40
1:K:36:ILE:HG22	1:K:37:LYS:N	2.84	0.40
1:6:37:LYS:O	1:6:38:LYS:C	2.59	0.40
1:C:115:ILE:HD13	1:C:115:ILE:HG21	2.18	0.40
6:E:2188:HOH:O	1:G:122:VAL:HG23	2.21	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	1	142/144 (99%)	140 (99%)	1 (1%)	1 (1%)	25	18
1	2	142/144 (99%)	139 (98%)	3 (2%)	0	100	100
1	3	140/144 (97%)	138 (99%)	2 (1%)	0	100	100
1	4	144/144 (100%)	142 (99%)	2 (1%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	5	141/144 (98%)	140 (99%)	1 (1%)	0	100	100
1	6	141/144 (98%)	139 (99%)	2 (1%)	0	100	100
1	7	140/144 (97%)	138 (99%)	2 (1%)	0	100	100
1	8	141/144 (98%)	139 (99%)	2 (1%)	0	100	100
1	9	141/144 (98%)	139 (99%)	2 (1%)	0	100	100
1	A	143/144 (99%)	141 (99%)	2 (1%)	0	100	100
1	B	142/144 (99%)	141 (99%)	1 (1%)	0	100	100
1	C	141/144 (98%)	140 (99%)	1 (1%)	0	100	100
1	D	141/144 (98%)	139 (99%)	2 (1%)	0	100	100
1	E	141/144 (98%)	139 (99%)	2 (1%)	0	100	100
1	F	143/144 (99%)	141 (99%)	2 (1%)	0	100	100
1	G	141/144 (98%)	138 (98%)	3 (2%)	0	100	100
1	H	140/144 (97%)	139 (99%)	1 (1%)	0	100	100
1	I	141/144 (98%)	139 (99%)	2 (1%)	0	100	100
1	J	141/144 (98%)	139 (99%)	2 (1%)	0	100	100
1	K	141/144 (98%)	140 (99%)	1 (1%)	0	100	100
1	L	140/144 (97%)	139 (99%)	1 (1%)	0	100	100
1	M	141/144 (98%)	139 (99%)	2 (1%)	0	100	100
1	N	140/144 (97%)	138 (99%)	2 (1%)	0	100	100
1	O	140/144 (97%)	139 (99%)	1 (1%)	0	100	100
1	P	141/144 (98%)	139 (99%)	2 (1%)	0	100	100
1	Q	142/144 (99%)	141 (99%)	1 (1%)	0	100	100
1	R	141/144 (98%)	139 (99%)	2 (1%)	0	100	100
1	S	142/144 (99%)	141 (99%)	1 (1%)	0	100	100
1	T	142/144 (99%)	140 (99%)	2 (1%)	0	100	100
1	U	142/144 (99%)	139 (98%)	3 (2%)	0	100	100
1	V	141/144 (98%)	139 (99%)	2 (1%)	0	100	100
1	W	140/144 (97%)	138 (99%)	2 (1%)	0	100	100
1	X	143/144 (99%)	142 (99%)	1 (1%)	0	100	100
1	Y	141/144 (98%)	140 (99%)	1 (1%)	0	100	100
1	Z	141/144 (98%)	139 (99%)	2 (1%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	a	142/144 (99%)	140 (99%)	2 (1%)	0	100	100
1	b	141/144 (98%)	139 (99%)	2 (1%)	0	100	100
1	c	142/144 (99%)	140 (99%)	2 (1%)	0	100	100
1	d	141/144 (98%)	139 (99%)	1 (1%)	1 (1%)	25	18
1	e	141/144 (98%)	139 (99%)	2 (1%)	0	100	100
1	f	141/144 (98%)	139 (99%)	2 (1%)	0	100	100
1	g	141/144 (98%)	139 (99%)	2 (1%)	0	100	100
1	h	140/144 (97%)	138 (99%)	2 (1%)	0	100	100
1	i	141/144 (98%)	137 (97%)	4 (3%)	0	100	100
1	j	142/144 (99%)	137 (96%)	5 (4%)	0	100	100
1	k	141/144 (98%)	136 (96%)	5 (4%)	0	100	100
1	l	142/144 (99%)	141 (99%)	1 (1%)	0	100	100
1	m	141/144 (98%)	139 (99%)	2 (1%)	0	100	100
All	All	6780/6912 (98%)	6685 (99%)	93 (1%)	2 (0%)	100	100

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	1	77	SER
1	d	77	SER

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	1	125/125 (100%)	122 (98%)	3 (2%)	54	56
1	2	125/125 (100%)	123 (98%)	2 (2%)	68	72
1	3	123/125 (98%)	120 (98%)	3 (2%)	54	56
1	4	127/125 (102%)	125 (98%)	2 (2%)	68	72
1	5	124/125 (99%)	122 (98%)	2 (2%)	68	72

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	6	124/125 (99%)	123 (99%)	1 (1%)	85	88
1	7	123/125 (98%)	121 (98%)	2 (2%)	68	72
1	8	124/125 (99%)	119 (96%)	5 (4%)	36	32
1	9	125/125 (100%)	120 (96%)	5 (4%)	36	32
1	A	126/125 (101%)	124 (98%)	2 (2%)	68	72
1	B	125/125 (100%)	123 (98%)	2 (2%)	68	72
1	C	125/125 (100%)	118 (94%)	7 (6%)	25	19
1	D	124/125 (99%)	120 (97%)	4 (3%)	44	42
1	E	124/125 (99%)	120 (97%)	4 (3%)	44	42
1	F	126/125 (101%)	123 (98%)	3 (2%)	54	56
1	G	125/125 (100%)	121 (97%)	4 (3%)	44	42
1	H	123/125 (98%)	117 (95%)	6 (5%)	29	24
1	I	124/125 (99%)	120 (97%)	4 (3%)	44	42
1	J	124/125 (99%)	119 (96%)	5 (4%)	36	32
1	K	124/125 (99%)	121 (98%)	3 (2%)	54	56
1	L	123/125 (98%)	121 (98%)	2 (2%)	68	72
1	M	124/125 (99%)	122 (98%)	2 (2%)	68	72
1	N	123/125 (98%)	121 (98%)	2 (2%)	68	72
1	O	123/125 (98%)	117 (95%)	6 (5%)	29	24
1	P	125/125 (100%)	122 (98%)	3 (2%)	54	56
1	Q	125/125 (100%)	122 (98%)	3 (2%)	54	56
1	R	124/125 (99%)	120 (97%)	4 (3%)	44	42
1	S	125/125 (100%)	118 (94%)	7 (6%)	25	19
1	T	125/125 (100%)	121 (97%)	4 (3%)	44	42
1	U	125/125 (100%)	123 (98%)	2 (2%)	68	72
1	V	124/125 (99%)	120 (97%)	4 (3%)	44	42
1	W	123/125 (98%)	120 (98%)	3 (2%)	54	56
1	X	126/125 (101%)	124 (98%)	2 (2%)	68	72
1	Y	124/125 (99%)	119 (96%)	5 (4%)	36	32
1	Z	125/125 (100%)	122 (98%)	3 (2%)	54	56
1	a	125/125 (100%)	121 (97%)	4 (3%)	44	42

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	b	124/125 (99%)	120 (97%)	4 (3%)	44	42
1	c	125/125 (100%)	122 (98%)	3 (2%)	54	56
1	d	124/125 (99%)	121 (98%)	3 (2%)	54	56
1	e	124/125 (99%)	119 (96%)	5 (4%)	36	32
1	f	124/125 (99%)	120 (97%)	4 (3%)	44	42
1	g	125/125 (100%)	120 (96%)	5 (4%)	36	32
1	h	123/125 (98%)	119 (97%)	4 (3%)	43	41
1	i	124/125 (99%)	121 (98%)	3 (2%)	54	56
1	j	125/125 (100%)	119 (95%)	6 (5%)	30	25
1	k	124/125 (99%)	117 (94%)	7 (6%)	25	19
1	l	125/125 (100%)	121 (97%)	4 (3%)	44	42
1	m	124/125 (99%)	121 (98%)	3 (2%)	54	56
All	All	5970/6000 (100%)	5794 (97%)	176 (3%)	48	47

All (176) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	1	12	GLU
1	1	35	THR
1	1	128	GLU
1	2	35	THR
1	2	103	GLU
1	3	12	GLU
1	3	35	THR
1	3	99	LEU
1	4	3	MET
1	4	12	GLU
1	5	12	GLU
1	5	35	THR
1	6	35	THR
1	7	35	THR
1	7	124	LEU
1	8	3	MET
1	8	12	GLU
1	8	35	THR
1	8	42	LYS
1	8	144	LYS
1	9	42	LYS

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Mol	Chain	Res	Type
1	9	124	LEU
1	9	129	HIS
1	9	144[A]	LYS
1	9	144[B]	LYS
1	A	35	THR
1	A	124	LEU
1	B	99	LEU
1	B	124	LEU
1	C	3	MET
1	C	18[A]	SER
1	C	18[B]	SER
1	C	35	THR
1	C	99	LEU
1	C	103	GLU
1	C	124	LEU
1	D	12	GLU
1	D	100	ARG
1	D	124	LEU
1	D	126	GLU
1	E	12	GLU
1	E	35	THR
1	E	99	LEU
1	E	124	LEU
1	F	12	GLU
1	F	35	THR
1	F	124	LEU
1	G	12	GLU
1	G	35	THR
1	G	103	GLU
1	G	124	LEU
1	H	12	GLU
1	H	35	THR
1	H	99	LEU
1	H	103	GLU
1	H	124	LEU
1	H	143	THR
1	I	103	GLU
1	I	119[A]	MET
1	I	119[B]	MET
1	I	124	LEU
1	J	12	GLU
1	J	35	THR

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Mol	Chain	Res	Type
1	J	54	GLU
1	J	103	GLU
1	J	124	LEU
1	K	12	GLU
1	K	100	ARG
1	K	124	LEU
1	L	12	GLU
1	L	124	LEU
1	M	12	GLU
1	M	124	LEU
1	N	12	GLU
1	N	124	LEU
1	O	12	GLU
1	O	18	SER
1	O	32	GLU
1	O	35	THR
1	O	99	LEU
1	O	124	LEU
1	P	12	GLU
1	P	35	THR
1	P	124	LEU
1	Q	12	GLU
1	Q	35	THR
1	Q	124	LEU
1	R	12	GLU
1	R	35	THR
1	R	105	LYS
1	R	124	LEU
1	S	12	GLU
1	S	35	THR
1	S	105	LYS
1	S	124	LEU
1	S	129[A]	HIS
1	S	129[B]	HIS
1	S	143	THR
1	T	12	GLU
1	T	35	THR
1	T	100	ARG
1	T	124	LEU
1	U	12	GLU
1	U	124	LEU
1	V	12	GLU

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Mol	Chain	Res	Type
1	V	35	THR
1	V	42	LYS
1	V	124	LEU
1	W	12	GLU
1	W	35	THR
1	W	124	LEU
1	X	12	GLU
1	X	124	LEU
1	Y	12	GLU
1	Y	35	THR
1	Y	100	ARG
1	Y	124	LEU
1	Y	143	THR
1	Z	35	THR
1	Z	99	LEU
1	Z	124	LEU
1	a	12	GLU
1	a	40	GLU
1	a	124	LEU
1	a	144	LYS
1	b	12	GLU
1	b	54	GLU
1	b	124	LEU
1	b	144	LYS
1	c	35	THR
1	c	124	LEU
1	c	144	LYS
1	d	3	MET
1	d	35	THR
1	d	124	LEU
1	e	12	GLU
1	e	35	THR
1	e	42	LYS
1	e	124	LEU
1	e	144	LYS
1	f	12	GLU
1	f	37	LYS
1	f	42	LYS
1	f	124	LEU
1	g	35	THR
1	g	42	LYS
1	g	124	LEU

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Mol	Chain	Res	Type
1	g	144[A]	LYS
1	g	144[B]	LYS
1	h	12	GLU
1	h	35	THR
1	h	124	LEU
1	h	144	LYS
1	i	35	THR
1	i	99	LEU
1	i	124	LEU
1	j	12	GLU
1	j	35	THR
1	j	42	LYS
1	j	54	GLU
1	j	124	LEU
1	j	144	LYS
1	k	3	MET
1	k	12	GLU
1	k	42	LYS
1	k	99	LEU
1	k	102	THR
1	k	124	LEU
1	k	144	LYS
1	l	35	THR
1	l	42	LYS
1	l	124	LEU
1	l	144	LYS
1	m	12	GLU
1	m	124	LEU
1	m	144	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (44) such sidechains are listed below:

Mol	Chain	Res	Type
1	1	4	GLN
1	2	112	GLN
1	3	4	GLN
1	4	4	GLN
1	4	112	GLN
1	5	4	GLN
1	7	4	GLN
1	8	4	GLN
1	9	4	GLN

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Mol	Chain	Res	Type
1	A	4	GLN
1	B	4	GLN
1	C	4	GLN
1	D	4	GLN
1	F	4	GLN
1	G	4	GLN
1	H	4	GLN
1	I	4	GLN
1	I	112	GLN
1	J	4	GLN
1	K	4	GLN
1	L	4	GLN
1	M	4	GLN
1	M	129	HIS
1	N	4	GLN
1	O	4	GLN
1	P	4	GLN
1	Q	4	GLN
1	S	4	GLN
1	U	4	GLN
1	V	4	GLN
1	W	4	GLN
1	X	4	GLN
1	Y	4	GLN
1	Z	4	GLN
1	a	4	GLN
1	b	4	GLN
1	d	4	GLN
1	e	4	GLN
1	f	4	GLN
1	g	4	GLN
1	h	4	GLN
1	k	4	GLN
1	l	4	GLN
1	m	4	GLN

5.3.3 RNA ⓘ

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 146 ligands modelled in this entry, 48 are monoatomic - leaving 98 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$
2	DUR	1	1145	-	13,17,17	1.28	2 (15%)	16,24,24	3.85	4 (25%)
3	POP	1	1146	4	8,8,8	0.81	0	8,13,13	1.39	2 (25%)
2	DUR	2	1145	-	13,17,17	0.87	1 (7%)	16,24,24	2.16	1 (6%)
3	POP	2	1146	4	8,8,8	1.65	1 (12%)	8,13,13	0.86	0
2	DUR	3	1145	-	13,17,17	1.02	1 (7%)	16,24,24	2.18	2 (12%)
3	POP	3	1146	4	8,8,8	1.85	1 (12%)	8,13,13	1.64	1 (12%)
2	DUR	4	1145	-	13,17,17	0.91	0	16,24,24	2.25	3 (18%)
3	POP	4	1146	4	8,8,8	1.75	2 (25%)	8,13,13	1.90	4 (50%)
2	DUR	5	1145	-	13,17,17	1.37	1 (7%)	16,24,24	2.12	2 (12%)
3	POP	5	1146	4	8,8,8	1.06	0	8,13,13	1.12	1 (12%)
2	DUR	6	1145	-	13,17,17	0.96	1 (7%)	16,24,24	2.64	1 (6%)
3	POP	6	1146	4	8,8,8	1.75	1 (12%)	8,13,13	1.42	1 (12%)
2	DUR	7	1145	-	13,17,17	1.18	2 (15%)	16,24,24	2.11	1 (6%)
3	POP	7	1146	4	8,8,8	0.76	0	8,13,13	1.12	0
2	DUR	8	1145	-	13,17,17	1.26	1 (7%)	16,24,24	2.83	1 (6%)
3	POP	8	1146	4	8,8,8	1.61	1 (12%)	8,13,13	1.40	1 (12%)
2	DUR	9	1145	-	13,17,17	0.87	0	16,24,24	2.25	2 (12%)
3	POP	9	1146	4	8,8,8	2.07	2 (25%)	8,13,13	1.21	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	DUR	A	1145	-	13,17,17	1.21	1 (7%)	16,24,24	3.06	5 (31%)
3	POP	A	1146	4	8,8,8	1.16	0	8,13,13	1.29	1 (12%)
2	DUR	B	1145	-	13,17,17	0.83	1 (7%)	16,24,24	2.27	1 (6%)
3	POP	B	1146	4	8,8,8	2.09	1 (12%)	8,13,13	1.84	2 (25%)
2	DUR	C	1145	-	13,17,17	1.15	2 (15%)	16,24,24	1.94	1 (6%)
3	POP	C	1146	4	8,8,8	0.58	0	8,13,13	1.62	2 (25%)
2	DUR	D	1145	-	13,17,17	1.06	1 (7%)	16,24,24	2.45	4 (25%)
3	POP	D	1146	4	8,8,8	1.42	1 (12%)	8,13,13	1.58	3 (37%)
2	DUR	E	1145	-	13,17,17	1.00	1 (7%)	16,24,24	2.50	4 (25%)
3	POP	E	1146	4	8,8,8	0.91	0	8,13,13	1.31	1 (12%)
2	DUR	F	1145	-	13,17,17	1.67	2 (15%)	16,24,24	3.26	3 (18%)
3	POP	F	1146	4	8,8,8	1.88	3 (37%)	8,13,13	1.53	1 (12%)
2	DUR	G	1145	-	13,17,17	1.02	1 (7%)	16,24,24	2.97	3 (18%)
3	POP	G	1146	4	8,8,8	2.09	1 (12%)	8,13,13	2.05	4 (50%)
2	DUR	H	1145	-	13,17,17	1.05	1 (7%)	16,24,24	2.59	4 (25%)
3	POP	H	1146	4	8,8,8	1.34	0	8,13,13	1.37	0
2	DUR	I	1145	-	13,17,17	1.83	2 (15%)	16,24,24	1.87	1 (6%)
3	POP	I	1146	4	8,8,8	1.52	2 (25%)	8,13,13	1.52	1 (12%)
2	DUR	J	1145	-	13,17,17	1.60	3 (23%)	16,24,24	2.75	4 (25%)
3	POP	J	1146	4	8,8,8	0.81	0	8,13,13	1.14	1 (12%)
2	DUR	K	1145	-	13,17,17	1.25	1 (7%)	16,24,24	1.72	1 (6%)
3	POP	K	1146	4	8,8,8	1.65	2 (25%)	8,13,13	1.11	0
2	DUR	L	1145	-	13,17,17	1.44	2 (15%)	16,24,24	2.20	3 (18%)
3	POP	L	1146	4	8,8,8	1.09	1 (12%)	8,13,13	1.18	0
2	DUR	M	1145	-	13,17,17	1.10	1 (7%)	16,24,24	1.54	1 (6%)
3	POP	M	1146	4	8,8,8	0.95	0	8,13,13	1.67	2 (25%)
2	DUR	N	1145	-	13,17,17	0.91	0	16,24,24	2.34	4 (25%)
3	POP	N	1146	4	8,8,8	1.82	1 (12%)	8,13,13	1.45	2 (25%)
2	DUR	O	1145	-	13,17,17	1.34	1 (7%)	16,24,24	2.86	3 (18%)
3	POP	O	1146	4	8,8,8	1.43	1 (12%)	8,13,13	1.46	2 (25%)
2	DUR	P	1145	-	13,17,17	0.65	0	16,24,24	2.23	2 (12%)
3	POP	P	1146	4	8,8,8	1.43	1 (12%)	8,13,13	1.87	3 (37%)
2	DUR	Q	1145	-	13,17,17	0.97	0	16,24,24	1.73	1 (6%)
3	POP	Q	1146	4	8,8,8	1.37	2 (25%)	8,13,13	1.19	1 (12%)
2	DUR	R	1145	-	13,17,17	1.63	1 (7%)	16,24,24	2.26	1 (6%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	POP	R	1146	4	8,8,8	1.06	1 (12%)	8,13,13	1.76	3 (37%)
2	DUR	S	1145	-	13,17,17	1.27	1 (7%)	16,24,24	2.56	2 (12%)
3	POP	S	1146	4	8,8,8	2.06	1 (12%)	8,13,13	0.98	0
2	DUR	T	1145	-	13,17,17	0.92	1 (7%)	16,24,24	2.14	3 (18%)
3	POP	T	1146	4	8,8,8	1.09	1 (12%)	8,13,13	1.56	2 (25%)
2	DUR	U	1145	-	13,17,17	0.99	0	16,24,24	1.79	2 (12%)
3	POP	U	1146	4	8,8,8	0.90	0	8,13,13	1.33	0
2	DUR	V	1145	-	13,17,17	1.39	2 (15%)	16,24,24	2.09	2 (12%)
3	POP	V	1146	4	8,8,8	0.96	0	8,13,13	1.78	3 (37%)
2	DUR	W	1145	-	13,17,17	1.03	1 (7%)	16,24,24	2.59	1 (6%)
3	POP	W	1146	4	8,8,8	1.13	0	8,13,13	1.48	1 (12%)
2	DUR	X	1145	-	13,17,17	1.15	1 (7%)	16,24,24	2.50	2 (12%)
3	POP	X	1146	4	8,8,8	1.47	1 (12%)	8,13,13	1.67	2 (25%)
2	DUR	Y	1145	-	13,17,17	1.13	1 (7%)	16,24,24	2.69	3 (18%)
3	POP	Y	1146	4	8,8,8	0.95	0	8,13,13	1.50	2 (25%)
2	DUR	Z	1145	-	13,17,17	1.20	1 (7%)	16,24,24	1.95	2 (12%)
3	POP	Z	1146	4	8,8,8	0.91	0	8,13,13	1.41	2 (25%)
2	DUR	a	1145	-	13,17,17	1.12	1 (7%)	16,24,24	2.14	3 (18%)
3	POP	a	1146	4	8,8,8	1.04	1 (12%)	8,13,13	1.58	1 (12%)
2	DUR	b	1145	-	13,17,17	1.44	1 (7%)	16,24,24	2.86	3 (18%)
3	POP	b	1146	4	8,8,8	1.97	2 (25%)	8,13,13	1.51	2 (25%)
2	DUR	c	1145	-	13,17,17	1.30	2 (15%)	16,24,24	2.45	1 (6%)
3	POP	c	1146	4	8,8,8	0.65	0	8,13,13	1.99	3 (37%)
2	DUR	d	1145	-	13,17,17	1.00	1 (7%)	16,24,24	2.72	2 (12%)
3	POP	d	1146	4	8,8,8	0.95	0	8,13,13	1.90	4 (50%)
2	DUR	e	1145	-	13,17,17	1.44	1 (7%)	16,24,24	2.59	3 (18%)
3	POP	e	1146	4	8,8,8	1.59	2 (25%)	8,13,13	1.51	2 (25%)
5	PO4	e	1148	-	4,4,4	0.52	0	6,6,6	0.71	0
2	DUR	f	1145	-	13,17,17	0.83	0	16,24,24	2.52	2 (12%)
3	POP	f	1146	4	8,8,8	0.61	0	8,13,13	1.82	3 (37%)
2	DUR	g	1145	-	13,17,17	0.90	1 (7%)	16,24,24	2.54	2 (12%)
3	POP	g	1146	4	8,8,8	1.37	1 (12%)	8,13,13	1.31	1 (12%)
2	DUR	h	1145	-	13,17,17	1.81	2 (15%)	16,24,24	1.92	1 (6%)
3	POP	h	1146	4	8,8,8	1.27	2 (25%)	8,13,13	1.54	2 (25%)
2	DUR	i	1145	-	13,17,17	1.05	1 (7%)	16,24,24	2.20	2 (12%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	POP	i	1146	4	8,8,8	1.18	1 (12%)	8,13,13	1.31	1 (12%)
2	DUR	j	1145	-	13,17,17	1.16	1 (7%)	16,24,24	2.40	3 (18%)
3	POP	j	1146	4	8,8,8	1.65	2 (25%)	8,13,13	1.96	2 (25%)
2	DUR	k	1145	-	13,17,17	0.92	1 (7%)	16,24,24	2.15	3 (18%)
3	POP	k	1146	4	8,8,8	1.15	1 (12%)	8,13,13	1.46	1 (12%)
2	DUR	l	1145	-	13,17,17	1.46	3 (23%)	16,24,24	2.46	1 (6%)
3	POP	l	1146	4	8,8,8	1.31	1 (12%)	8,13,13	1.14	1 (12%)
5	PO4	l	1148	-	4,4,4	1.37	1 (25%)	6,6,6	0.98	0
2	DUR	m	1145	-	13,17,17	0.99	1 (7%)	16,24,24	2.75	1 (6%)
3	POP	m	1146	4	8,8,8	1.51	1 (12%)	8,13,13	1.74	1 (12%)

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
2	DUR	1	1145	-	-	0/2/18/18	0/2/2/2
3	POP	1	1146	4	-	0/6/6/6	0/0/0/0
2	DUR	2	1145	-	-	0/2/18/18	0/2/2/2
3	POP	2	1146	4	-	0/6/6/6	0/0/0/0
2	DUR	3	1145	-	-	0/2/18/18	0/2/2/2
3	POP	3	1146	4	-	0/6/6/6	0/0/0/0
2	DUR	4	1145	-	-	0/2/18/18	0/2/2/2
3	POP	4	1146	4	-	0/6/6/6	0/0/0/0
2	DUR	5	1145	-	-	0/2/18/18	0/2/2/2
3	POP	5	1146	4	-	0/6/6/6	0/0/0/0
2	DUR	6	1145	-	-	0/2/18/18	0/2/2/2
3	POP	6	1146	4	-	0/6/6/6	0/0/0/0
2	DUR	7	1145	-	-	0/2/18/18	0/2/2/2
3	POP	7	1146	4	-	0/6/6/6	0/0/0/0
2	DUR	8	1145	-	-	0/2/18/18	0/2/2/2
3	POP	8	1146	4	-	0/6/6/6	0/0/0/0
2	DUR	9	1145	-	-	0/2/18/18	0/2/2/2
3	POP	9	1146	4	-	0/6/6/6	0/0/0/0
2	DUR	A	1145	-	-	0/2/18/18	0/2/2/2
3	POP	A	1146	4	-	0/6/6/6	0/0/0/0
2	DUR	B	1145	-	-	0/2/18/18	0/2/2/2
3	POP	B	1146	4	-	0/6/6/6	0/0/0/0
2	DUR	C	1145	-	-	0/2/18/18	0/2/2/2

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	POP	C	1146	4	-	0/6/6/6	0/0/0/0
2	DUR	D	1145	-	-	0/2/18/18	0/2/2/2
3	POP	D	1146	4	-	0/6/6/6	0/0/0/0
2	DUR	E	1145	-	-	0/2/18/18	0/2/2/2
3	POP	E	1146	4	-	0/6/6/6	0/0/0/0
2	DUR	F	1145	-	-	0/2/18/18	0/2/2/2
3	POP	F	1146	4	-	0/6/6/6	0/0/0/0
2	DUR	G	1145	-	-	0/2/18/18	0/2/2/2
3	POP	G	1146	4	-	0/6/6/6	0/0/0/0
2	DUR	H	1145	-	-	0/2/18/18	0/2/2/2
3	POP	H	1146	4	-	0/6/6/6	0/0/0/0
2	DUR	I	1145	-	-	0/2/18/18	0/2/2/2
3	POP	I	1146	4	-	0/6/6/6	0/0/0/0
2	DUR	J	1145	-	1/1/3/3	0/2/18/18	0/2/2/2
3	POP	J	1146	4	-	0/6/6/6	0/0/0/0
2	DUR	K	1145	-	-	0/2/18/18	0/2/2/2
3	POP	K	1146	4	-	0/6/6/6	0/0/0/0
2	DUR	L	1145	-	-	0/2/18/18	0/2/2/2
3	POP	L	1146	4	-	0/6/6/6	0/0/0/0
2	DUR	M	1145	-	-	0/2/18/18	0/2/2/2
3	POP	M	1146	4	-	0/6/6/6	0/0/0/0
2	DUR	N	1145	-	-	0/2/18/18	0/2/2/2
3	POP	N	1146	4	-	0/6/6/6	0/0/0/0
2	DUR	O	1145	-	-	0/2/18/18	0/2/2/2
3	POP	O	1146	4	-	0/6/6/6	0/0/0/0
2	DUR	P	1145	-	-	0/2/18/18	0/2/2/2
3	POP	P	1146	4	-	0/6/6/6	0/0/0/0
2	DUR	Q	1145	-	-	0/2/18/18	0/2/2/2
3	POP	Q	1146	4	-	0/6/6/6	0/0/0/0
2	DUR	R	1145	-	-	0/2/18/18	0/2/2/2
3	POP	R	1146	4	-	0/6/6/6	0/0/0/0
2	DUR	S	1145	-	-	0/2/18/18	0/2/2/2
3	POP	S	1146	4	-	0/6/6/6	0/0/0/0
2	DUR	T	1145	-	-	0/2/18/18	0/2/2/2
3	POP	T	1146	4	-	0/6/6/6	0/0/0/0
2	DUR	U	1145	-	-	0/2/18/18	0/2/2/2
3	POP	U	1146	4	-	0/6/6/6	0/0/0/0
2	DUR	V	1145	-	-	0/2/18/18	0/2/2/2
3	POP	V	1146	4	-	0/6/6/6	0/0/0/0
2	DUR	W	1145	-	-	0/2/18/18	0/2/2/2
3	POP	W	1146	4	-	0/6/6/6	0/0/0/0
2	DUR	X	1145	-	-	0/2/18/18	0/2/2/2

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	POP	X	1146	4	-	0/6/6/6	0/0/0/0
2	DUR	Y	1145	-	-	0/2/18/18	0/2/2/2
3	POP	Y	1146	4	-	0/6/6/6	0/0/0/0
2	DUR	Z	1145	-	-	0/2/18/18	0/2/2/2
3	POP	Z	1146	4	-	0/6/6/6	0/0/0/0
2	DUR	a	1145	-	-	0/2/18/18	0/2/2/2
3	POP	a	1146	4	-	0/6/6/6	0/0/0/0
2	DUR	b	1145	-	-	0/2/18/18	0/2/2/2
3	POP	b	1146	4	-	0/6/6/6	0/0/0/0
2	DUR	c	1145	-	-	0/2/18/18	0/2/2/2
3	POP	c	1146	4	-	0/6/6/6	0/0/0/0
2	DUR	d	1145	-	-	0/2/18/18	0/2/2/2
3	POP	d	1146	4	-	0/6/6/6	0/0/0/0
2	DUR	e	1145	-	-	0/2/18/18	0/2/2/2
3	POP	e	1146	4	-	0/6/6/6	0/0/0/0
5	PO4	e	1148	-	-	0/0/0/0	0/0/0/0
2	DUR	f	1145	-	-	0/2/18/18	0/2/2/2
3	POP	f	1146	4	-	0/6/6/6	0/0/0/0
2	DUR	g	1145	-	-	0/2/18/18	0/2/2/2
3	POP	g	1146	4	-	0/6/6/6	0/0/0/0
2	DUR	h	1145	-	-	0/2/18/18	0/2/2/2
3	POP	h	1146	4	-	0/6/6/6	0/0/0/0
2	DUR	i	1145	-	-	0/2/18/18	0/2/2/2
3	POP	i	1146	4	-	0/6/6/6	0/0/0/0
2	DUR	j	1145	-	-	0/2/18/18	0/2/2/2
3	POP	j	1146	4	-	0/6/6/6	0/0/0/0
2	DUR	k	1145	-	-	0/2/18/18	0/2/2/2
3	POP	k	1146	4	-	0/6/6/6	0/0/0/0
2	DUR	l	1145	-	-	0/2/18/18	0/2/2/2
3	POP	l	1146	4	-	0/6/6/6	0/0/0/0
5	PO4	l	1148	-	-	0/0/0/0	0/0/0/0
2	DUR	m	1145	-	-	0/2/18/18	0/2/2/2
3	POP	m	1146	4	-	0/6/6/6	0/0/0/0

All (97) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	h	1145	DUR	C2-N3	-5.15	1.28	1.38
2	R	1145	DUR	C6-N1	-5.01	1.29	1.35
2	F	1145	DUR	C2-N3	-4.66	1.28	1.38
2	I	1145	DUR	C2-N3	-4.56	1.29	1.38
2	J	1145	DUR	C2-N3	-4.51	1.29	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	e	1145	DUR	C2-N3	-4.18	1.29	1.38
2	5	1145	DUR	C6-N1	-4.05	1.30	1.35
2	b	1145	DUR	C2-N3	-4.02	1.30	1.38
2	c	1145	DUR	C6-N1	-3.99	1.30	1.35
2	8	1145	DUR	C2-N3	-3.89	1.30	1.38
2	O	1145	DUR	C2-N3	-3.89	1.30	1.38
2	I	1145	DUR	C6-N1	-3.86	1.30	1.35
2	L	1145	DUR	C2-N3	-3.83	1.30	1.38
2	Z	1145	DUR	C2-N3	-3.56	1.31	1.38
2	S	1145	DUR	C2-N3	-3.42	1.31	1.38
2	V	1145	DUR	C6-N1	-3.31	1.31	1.35
2	X	1145	DUR	C2-N3	-3.28	1.31	1.38
2	Y	1145	DUR	C2-N3	-3.26	1.31	1.38
2	V	1145	DUR	C2-N3	-3.17	1.31	1.38
2	E	1145	DUR	C2-N3	-3.12	1.32	1.38
2	A	1145	DUR	C2-N3	-3.04	1.32	1.38
2	3	1145	DUR	C2-N3	-2.87	1.32	1.38
2	l	1145	DUR	C2-N3	-2.84	1.32	1.38
2	W	1145	DUR	C2-N3	-2.84	1.32	1.38
2	M	1145	DUR	C2-N3	-2.81	1.32	1.38
2	j	1145	DUR	C2-N3	-2.75	1.32	1.38
2	2	1145	DUR	C2-N3	-2.63	1.33	1.38
2	7	1145	DUR	C6-N1	-2.58	1.32	1.35
2	6	1145	DUR	C2-N3	-2.54	1.33	1.38
2	a	1145	DUR	C2-N3	-2.51	1.33	1.38
2	H	1145	DUR	C2-N3	-2.46	1.33	1.38
3	R	1146	POP	P2-O	-2.45	1.56	1.60
2	d	1145	DUR	C2-N3	-2.38	1.33	1.38
2	K	1145	DUR	C2-N3	-2.34	1.33	1.38
2	L	1145	DUR	C6-N1	-2.31	1.32	1.35
2	T	1145	DUR	C2-N3	-2.30	1.33	1.38
2	7	1145	DUR	C2-N3	-2.28	1.33	1.38
2	J	1145	DUR	C2'-C3'	-2.27	1.46	1.52
2	h	1145	DUR	C6-N1	-2.27	1.32	1.35
2	l	1145	DUR	O4'-C4'	-2.25	1.39	1.45
2	m	1145	DUR	C2-N3	-2.22	1.33	1.38
3	F	1146	POP	P2-O4	-2.19	1.43	1.50
3	I	1146	POP	P2-O6	-2.17	1.45	1.54
2	B	1145	DUR	C2-N3	-2.15	1.33	1.38
3	h	1146	POP	P2-O5	-2.15	1.46	1.54
2	k	1145	DUR	C2-N3	-2.14	1.33	1.38
2	g	1145	DUR	C2-N3	-2.13	1.33	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	1145	DUR	C6-N1	-2.13	1.32	1.35
2	c	1145	DUR	C2-N3	-2.10	1.34	1.38
3	e	1146	POP	P2-O6	-2.06	1.46	1.54
2	i	1145	DUR	C2-N3	-2.05	1.34	1.38
2	C	1145	DUR	C2-N3	-2.04	1.34	1.38
3	T	1146	POP	P1-O	2.01	1.63	1.60
2	F	1145	DUR	O3'-C3'	2.02	1.47	1.43
3	Q	1146	POP	P2-O	2.03	1.63	1.60
2	J	1145	DUR	C4-N3	2.07	1.36	1.33
3	K	1146	POP	P2-O	2.10	1.63	1.60
3	L	1146	POP	P1-O	2.10	1.63	1.60
3	b	1146	POP	P2-O	2.11	1.63	1.60
3	h	1146	POP	P1-O	2.14	1.63	1.60
5	l	1148	PO4	P-O1	2.23	1.55	1.50
3	a	1146	POP	P1-O	2.24	1.63	1.60
3	Q	1146	POP	P1-O	2.25	1.63	1.60
3	i	1146	POP	P2-O	2.34	1.63	1.60
2	l	1145	DUR	O4-C4	2.49	1.30	1.24
3	k	1146	POP	P1-O	2.50	1.64	1.60
3	I	1146	POP	P1-O	2.53	1.64	1.60
3	F	1146	POP	P1-O	2.56	1.64	1.60
2	G	1145	DUR	O4-C4	2.57	1.31	1.24
2	D	1145	DUR	C4-N3	2.58	1.37	1.33
2	l	1145	DUR	C6-C5	2.60	1.43	1.38
3	j	1146	POP	P1-O	2.68	1.64	1.60
3	K	1146	POP	P1-O	2.69	1.64	1.60
3	D	1146	POP	P2-O	2.69	1.64	1.60
3	e	1146	POP	P1-O	2.82	1.64	1.60
3	4	1146	POP	P2-O	2.83	1.64	1.60
3	l	1146	POP	P1-O	2.91	1.64	1.60
2	l	1145	DUR	O4'-C1'	2.94	1.49	1.42
3	P	1146	POP	P2-O	3.02	1.64	1.60
3	X	1146	POP	P2-O	3.04	1.65	1.60
3	O	1146	POP	P1-O	3.07	1.65	1.60
3	4	1146	POP	P1-O	3.09	1.65	1.60
3	g	1146	POP	P2-O	3.25	1.65	1.60
3	j	1146	POP	P2-O	3.42	1.65	1.60
3	9	1146	POP	P2-O	3.72	1.66	1.60
3	8	1146	POP	P2-O	3.73	1.66	1.60
3	m	1146	POP	P2-O	3.74	1.66	1.60
3	2	1146	POP	P2-O	3.91	1.66	1.60
3	9	1146	POP	P1-O	3.97	1.66	1.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	F	1146	POP	P2-O	4.01	1.66	1.60
3	6	1146	POP	P1-O	4.18	1.66	1.60
3	3	1146	POP	P1-O	4.22	1.66	1.60
3	N	1146	POP	P1-O	4.38	1.67	1.60
3	b	1146	POP	P1-O	4.50	1.67	1.60
3	S	1146	POP	P1-O	4.61	1.67	1.60
3	G	1146	POP	P2-O	4.98	1.68	1.60
3	B	1146	POP	P2-O	5.27	1.68	1.60

All (182) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1145	DUR	C6-N1-C2	-3.33	115.88	121.28
2	1	1145	DUR	C5-C4-N3	-2.92	116.16	123.12
2	k	1145	DUR	C2'-C1'-N1	-2.87	107.46	114.23
2	P	1145	DUR	C2'-C3'-C4'	-2.71	96.98	102.73
2	H	1145	DUR	C2'-C3'-C4'	-2.69	97.01	102.73
2	b	1145	DUR	O4'-C4'-C3'	-2.61	99.42	105.68
3	f	1146	POP	O-P1-O1	-2.59	95.53	111.44
2	H	1145	DUR	C6-N1-C2	-2.57	117.11	121.28
2	E	1145	DUR	C2'-C1'-N1	-2.54	108.23	114.23
2	b	1145	DUR	O4'-C1'-N1	-2.54	103.50	107.78
2	X	1145	DUR	O4'-C4'-C5'	-2.50	103.82	109.16
2	1	1145	DUR	C6-N1-C2	-2.47	117.28	121.28
2	V	1145	DUR	C6-N1-C2	-2.46	117.30	121.28
2	D	1145	DUR	C6-N1-C2	-2.45	117.31	121.28
3	I	1146	POP	O3-P1-O1	-2.42	101.05	110.50
2	9	1145	DUR	C5-C4-N3	-2.40	117.38	123.12
2	E	1145	DUR	O4'-C1'-N1	-2.39	103.75	107.78
2	O	1145	DUR	C6-N1-C2	-2.39	117.41	121.28
3	X	1146	POP	O-P1-O1	-2.35	97.00	111.44
2	j	1145	DUR	C6-N1-C2	-2.33	117.50	121.28
3	h	1146	POP	O-P1-O1	-2.28	97.42	111.44
2	Z	1145	DUR	C5'-C4'-C3'	-2.24	109.16	114.84
3	V	1146	POP	O6-P2-O4	-2.24	101.75	110.50
2	g	1145	DUR	C6-N1-C2	-2.23	117.67	121.28
2	e	1145	DUR	C6-N1-C2	-2.21	117.70	121.28
2	G	1145	DUR	C5-C4-N3	-2.19	117.89	123.12
2	T	1145	DUR	O4'-C4'-C5'	-2.17	104.51	109.16
3	R	1146	POP	O-P1-O1	-2.16	98.16	111.44
3	M	1146	POP	O-P1-O1	-2.16	98.16	111.44
2	A	1145	DUR	C5-C4-N3	-2.14	118.00	123.12

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	L	1145	DUR	C2'-C3'-C4'	-2.14	98.18	102.73
3	b	1146	POP	O-P1-O1	-2.14	98.28	111.44
3	D	1146	POP	O-P1-O1	-2.14	98.31	111.44
3	4	1146	POP	O-P1-O1	-2.13	98.32	111.44
2	a	1145	DUR	C6-N1-C2	-2.13	117.83	121.28
2	3	1145	DUR	O4'-C1'-N1	-2.12	104.21	107.78
3	N	1146	POP	O-P1-O1	-2.12	98.39	111.44
2	d	1145	DUR	C5-C4-N3	-2.12	118.07	123.12
3	W	1146	POP	O-P2-O4	-2.11	98.48	111.44
2	A	1145	DUR	C2'-C1'-N1	-2.10	109.27	114.23
2	5	1145	DUR	C5-C4-N3	-2.10	118.10	123.12
3	G	1146	POP	O-P1-O1	-2.10	98.53	111.44
2	F	1145	DUR	C5-C4-N3	-2.08	118.16	123.12
2	Y	1145	DUR	C6-N1-C2	-2.08	117.91	121.28
2	D	1145	DUR	C5-C4-N3	-2.06	118.19	123.12
3	e	1146	POP	O6-P2-O4	-2.04	102.50	110.50
3	Y	1146	POP	O-P2-O4	-2.03	98.94	111.44
2	N	1145	DUR	C5-C4-N3	-2.03	118.28	123.12
2	4	1145	DUR	C5-C4-N3	-2.02	118.29	123.12
2	j	1145	DUR	C2'-C1'-N1	-2.02	109.47	114.23
3	c	1146	POP	O-P1-O1	-2.01	99.05	111.44
3	d	1146	POP	O6-P2-O5	2.00	115.70	107.61
2	F	1145	DUR	O3'-C3'-C2'	2.01	118.14	110.83
2	1	1145	DUR	C5'-C4'-C3'	2.02	119.96	114.84
2	G	1145	DUR	O4'-C1'-N1	2.02	111.19	107.78
3	5	1146	POP	O3-P1-O2	2.03	115.78	107.61
3	V	1146	POP	O5-P2-O4	2.03	118.45	110.50
3	4	1146	POP	O5-P2-O4	2.04	118.47	110.50
3	h	1146	POP	O6-P2-O5	2.04	115.85	107.61
2	4	1145	DUR	O3'-C3'-C4'	2.05	118.24	110.13
3	J	1146	POP	O6-P2-O5	2.07	115.97	107.61
3	4	1146	POP	O6-P2-O5	2.07	115.98	107.61
3	A	1146	POP	O6-P2-O5	2.08	116.02	107.61
3	P	1146	POP	O6-P2-O5	2.09	116.06	107.61
2	A	1145	DUR	O4'-C1'-N1	2.11	111.34	107.78
3	C	1146	POP	O6-P2-O5	2.13	116.22	107.61
3	l	1146	POP	O3-P1-O1	2.14	118.89	110.50
3	Z	1146	POP	O3-P1-O1	2.16	118.97	110.50
3	D	1146	POP	O2-P1-O1	2.17	118.98	110.50
3	d	1146	POP	O3-P1-O1	2.18	119.02	110.50
3	C	1146	POP	O3-P1-O1	2.19	119.07	110.50
3	8	1146	POP	O6-P2-O5	2.20	116.50	107.61

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	N	1145	DUR	O4'-C1'-N1	2.21	111.50	107.78
2	Y	1145	DUR	O4'-C4'-C5'	2.22	113.91	109.16
3	1	1146	POP	O6-P2-O5	2.22	116.56	107.61
2	L	1145	DUR	C3'-C2'-C1'	2.22	108.17	102.48
3	e	1146	POP	O3-P1-O1	2.22	119.20	110.50
2	N	1145	DUR	C4'-O4'-C1'	2.23	114.87	109.42
3	G	1146	POP	O3-P1-O1	2.24	119.25	110.50
3	G	1146	POP	O6-P2-O5	2.24	116.63	107.61
3	O	1146	POP	O6-P2-O5	2.24	116.66	107.61
3	Q	1146	POP	O2-P1-O1	2.24	119.28	110.50
3	Z	1146	POP	O5-P2-O4	2.25	119.29	110.50
3	g	1146	POP	O6-P2-O5	2.25	116.69	107.61
2	a	1145	DUR	O3'-C3'-C4'	2.26	119.05	110.13
3	b	1146	POP	O6-P2-O5	2.30	116.89	107.61
2	e	1145	DUR	C3'-C2'-C1'	2.30	108.38	102.48
2	D	1145	DUR	O4'-C1'-N1	2.31	111.68	107.78
3	f	1146	POP	O5-P2-O4	2.32	119.58	110.50
3	6	1146	POP	O3-P1-O1	2.32	119.60	110.50
3	T	1146	POP	O6-P2-O5	2.33	117.03	107.61
3	Y	1146	POP	O6-P2-O5	2.35	117.10	107.61
3	1	1146	POP	O3-P1-O2	2.36	117.14	107.61
2	O	1145	DUR	O4'-C1'-N1	2.38	111.79	107.78
3	d	1146	POP	O3-P1-O2	2.38	117.22	107.61
3	d	1146	POP	O5-P2-O4	2.39	119.85	110.50
3	N	1146	POP	O3-P1-O1	2.40	119.91	110.50
2	E	1145	DUR	C3'-C2'-C1'	2.44	108.73	102.48
2	H	1145	DUR	O3'-C3'-C4'	2.45	119.80	110.13
2	U	1145	DUR	O3'-C3'-C2'	2.45	119.75	110.83
3	P	1146	POP	O5-P2-O4	2.47	120.15	110.50
2	k	1145	DUR	O4'-C1'-C2'	2.52	111.08	106.25
3	R	1146	POP	O3-P1-O2	2.52	117.79	107.61
3	k	1146	POP	O5-P2-O4	2.53	120.42	110.50
3	F	1146	POP	O3-P1-O1	2.53	120.42	110.50
3	B	1146	POP	O6-P2-O5	2.57	117.97	107.61
3	B	1146	POP	O5-P2-O4	2.59	120.65	110.50
3	E	1146	POP	O6-P2-O5	2.61	118.15	107.61
2	T	1145	DUR	C5'-C4'-C3'	2.65	121.58	114.84
2	f	1145	DUR	O3'-C3'-C2'	2.70	120.65	110.83
3	X	1146	POP	O3-P1-O1	2.71	121.08	110.50
2	J	1145	DUR	O4'-C4'-C5'	2.71	114.97	109.16
3	O	1146	POP	O2-P1-O1	2.74	121.23	110.50
3	3	1146	POP	O3-P1-O1	2.76	121.29	110.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	i	1146	POP	O6-P2-O5	2.76	118.75	107.61
3	P	1146	POP	O3-P1-O1	2.80	121.46	110.50
2	S	1145	DUR	O3'-C3'-C4'	2.87	121.44	110.13
3	T	1146	POP	O3-P1-O2	2.89	119.28	107.61
3	f	1146	POP	O2-P1-O1	2.91	121.89	110.50
3	D	1146	POP	O6-P2-O5	2.94	119.46	107.61
3	j	1146	POP	O3-P1-O1	3.02	122.31	110.50
3	c	1146	POP	O5-P2-O4	3.02	122.32	110.50
2	J	1145	DUR	C2'-C3'-C4'	3.03	109.18	102.73
3	R	1146	POP	O2-P1-O1	3.09	122.60	110.50
3	a	1146	POP	O5-P2-O4	3.12	122.72	110.50
3	M	1146	POP	O5-P2-O4	3.15	122.84	110.50
3	4	1146	POP	O3-P1-O1	3.20	123.02	110.50
3	G	1146	POP	O5-P2-O4	3.28	123.34	110.50
3	c	1146	POP	O2-P1-O1	3.37	123.70	110.50
3	j	1146	POP	O5-P2-O4	3.41	123.85	110.50
2	i	1145	DUR	O4'-C1'-N1	3.54	113.74	107.78
2	J	1145	DUR	O3'-C3'-C4'	3.77	125.03	110.13
3	V	1146	POP	O6-P2-O5	3.83	123.06	107.61
3	m	1146	POP	O6-P2-O5	3.86	123.17	107.61
2	U	1145	DUR	C4-N3-C2	5.17	118.57	114.13
2	M	1145	DUR	C4-N3-C2	5.40	118.77	114.13
2	C	1145	DUR	C4-N3-C2	5.91	119.21	114.13
2	Q	1145	DUR	C4-N3-C2	6.06	119.34	114.13
2	Z	1145	DUR	C4-N3-C2	6.07	119.34	114.13
2	K	1145	DUR	C4-N3-C2	6.16	119.42	114.13
2	V	1145	DUR	C4-N3-C2	6.55	119.76	114.13
2	k	1145	DUR	C4-N3-C2	6.77	119.94	114.13
2	h	1145	DUR	C4-N3-C2	6.82	119.99	114.13
2	i	1145	DUR	C4-N3-C2	6.87	120.03	114.13
2	I	1145	DUR	C4-N3-C2	6.96	120.11	114.13
2	a	1145	DUR	C4-N3-C2	7.00	120.14	114.13
2	T	1145	DUR	C4-N3-C2	7.09	120.22	114.13
2	7	1145	DUR	C4-N3-C2	7.23	120.34	114.13
2	5	1145	DUR	C4-N3-C2	7.39	120.47	114.13
2	L	1145	DUR	C4-N3-C2	7.69	120.74	114.13
2	P	1145	DUR	C4-N3-C2	7.71	120.76	114.13
2	3	1145	DUR	C4-N3-C2	7.77	120.80	114.13
2	4	1145	DUR	C4-N3-C2	7.84	120.86	114.13
2	N	1145	DUR	C4-N3-C2	7.84	120.87	114.13
2	2	1145	DUR	C4-N3-C2	7.85	120.87	114.13
2	B	1145	DUR	C4-N3-C2	7.85	120.88	114.13

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	9	1145	DUR	C4-N3-C2	7.96	120.97	114.13
2	j	1145	DUR	C4-N3-C2	7.98	120.98	114.13
2	R	1145	DUR	C4-N3-C2	8.17	121.15	114.13
2	E	1145	DUR	C4-N3-C2	8.35	121.30	114.13
2	D	1145	DUR	C4-N3-C2	8.39	121.34	114.13
2	H	1145	DUR	C4-N3-C2	8.70	121.60	114.13
2	J	1145	DUR	C4-N3-C2	8.75	121.65	114.13
2	f	1145	DUR	C4-N3-C2	8.80	121.69	114.13
2	X	1145	DUR	C4-N3-C2	8.85	121.73	114.13
2	c	1145	DUR	C4-N3-C2	9.02	121.88	114.13
2	l	1145	DUR	C4-N3-C2	9.09	121.94	114.13
2	S	1145	DUR	C4-N3-C2	9.13	121.97	114.13
2	e	1145	DUR	C4-N3-C2	9.29	122.11	114.13
2	g	1145	DUR	C4-N3-C2	9.34	122.16	114.13
2	Y	1145	DUR	C4-N3-C2	9.37	122.18	114.13
2	6	1145	DUR	C4-N3-C2	9.58	122.36	114.13
2	W	1145	DUR	C4-N3-C2	9.75	122.50	114.13
2	d	1145	DUR	C4-N3-C2	10.13	122.83	114.13
2	b	1145	DUR	C4-N3-C2	10.28	122.96	114.13
2	O	1145	DUR	C4-N3-C2	10.45	123.11	114.13
2	A	1145	DUR	C4-N3-C2	10.72	123.34	114.13
2	m	1145	DUR	C4-N3-C2	10.73	123.35	114.13
2	8	1145	DUR	C4-N3-C2	10.87	123.47	114.13
2	G	1145	DUR	C4-N3-C2	11.18	123.73	114.13
2	F	1145	DUR	C4-N3-C2	12.35	124.74	114.13
2	1	1145	DUR	C4-N3-C2	14.54	126.62	114.13

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	J	1145	DUR	C3'

There are no torsion outliers.

There are no ring outliers.

26 monomers are involved in 43 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	2	1145	DUR	3	0
2	4	1145	DUR	1	0
2	5	1145	DUR	1	0
2	7	1145	DUR	2	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	8	1145	DUR	2	0
2	A	1145	DUR	1	0
2	B	1145	DUR	2	0
2	C	1145	DUR	2	0
2	D	1145	DUR	1	0
2	E	1145	DUR	1	0
2	F	1145	DUR	2	0
2	G	1145	DUR	2	0
2	H	1145	DUR	2	0
2	I	1145	DUR	2	0
2	J	1145	DUR	2	0
2	L	1145	DUR	1	0
2	M	1145	DUR	2	0
2	N	1145	DUR	1	0
2	O	1145	DUR	2	0
2	P	1145	DUR	1	0
2	Q	1145	DUR	2	0
2	R	1145	DUR	1	0
2	T	1145	DUR	2	0
2	U	1145	DUR	1	0
2	X	1145	DUR	2	0
2	Y	1145	DUR	2	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	1	142/144 (98%)	-0.24	1 (0%) 87 87	14, 24, 44, 53	3 (2%)
1	2	142/144 (98%)	-0.45	0 100 100	13, 23, 39, 50	0
1	3	142/144 (98%)	-0.38	0 100 100	11, 21, 36, 44	1 (0%)
1	4	142/144 (98%)	-0.41	0 100 100	13, 21, 35, 49	0
1	5	142/144 (98%)	-0.21	1 (0%) 87 87	16, 27, 48, 56	2 (1%)
1	6	142/144 (98%)	-0.20	0 100 100	15, 26, 46, 58	3 (2%)
1	7	142/144 (98%)	-0.21	0 100 100	16, 27, 47, 60	0
1	8	142/144 (98%)	-0.44	0 100 100	13, 21, 40, 53	0
1	9	142/144 (98%)	-0.39	0 100 100	11, 22, 38, 53	1 (0%)
1	A	143/144 (99%)	-0.25	1 (0%) 87 87	16, 26, 44, 55	2 (1%)
1	B	142/144 (98%)	-0.33	0 100 100	14, 25, 43, 51	2 (1%)
1	C	142/144 (98%)	-0.16	2 (1%) 75 75	15, 25, 45, 57	2 (1%)
1	D	142/144 (98%)	-0.43	0 100 100	11, 22, 38, 54	2 (1%)
1	E	142/144 (98%)	-0.36	0 100 100	12, 22, 38, 57	1 (0%)
1	F	143/144 (99%)	-0.38	1 (0%) 87 87	11, 21, 40, 48	1 (0%)
1	G	142/144 (98%)	-0.39	0 100 100	11, 20, 37, 50	2 (1%)
1	H	142/144 (98%)	-0.42	1 (0%) 87 87	11, 20, 38, 48	1 (0%)
1	I	142/144 (98%)	-0.47	0 100 100	12, 20, 35, 47	2 (1%)
1	J	142/144 (98%)	-0.39	0 100 100	13, 23, 42, 51	1 (0%)
1	K	142/144 (98%)	-0.30	0 100 100	13, 24, 41, 52	4 (2%)
1	L	142/144 (98%)	-0.40	0 100 100	14, 23, 39, 55	0
1	M	142/144 (98%)	-0.34	1 (0%) 87 87	13, 23, 40, 48	1 (0%)
1	N	142/144 (98%)	-0.36	0 100 100	15, 25, 44, 52	1 (0%)
1	O	142/144 (98%)	-0.31	1 (0%) 87 87	16, 24, 41, 57	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	P	142/144 (98%)	-0.43	0 100 100	11, 20, 38, 48	1 (0%)
1	Q	143/144 (99%)	-0.37	0 100 100	13, 22, 40, 46	0
1	R	142/144 (98%)	-0.40	0 100 100	11, 22, 40, 50	1 (0%)
1	S	142/144 (98%)	-0.37	0 100 100	13, 23, 39, 54	0
1	T	142/144 (98%)	-0.38	0 100 100	13, 24, 45, 56	1 (0%)
1	U	142/144 (98%)	-0.38	0 100 100	13, 23, 42, 50	0
1	V	142/144 (98%)	-0.36	0 100 100	13, 23, 41, 50	1 (0%)
1	W	142/144 (98%)	-0.35	0 100 100	11, 22, 41, 51	0
1	X	142/144 (98%)	-0.42	0 100 100	12, 21, 41, 50	0
1	Y	142/144 (98%)	-0.26	0 100 100	14, 24, 44, 54	1 (0%)
1	Z	142/144 (98%)	-0.29	2 (1%) 75 75	14, 26, 43, 66	2 (1%)
1	a	143/144 (99%)	-0.43	0 100 100	13, 22, 36, 44	2 (1%)
1	b	143/144 (99%)	-0.37	0 100 100	12, 24, 45, 51	1 (0%)
1	c	142/144 (98%)	-0.31	0 100 100	13, 25, 44, 58	2 (1%)
1	d	142/144 (98%)	-0.30	1 (0%) 87 87	14, 25, 45, 60	0
1	e	142/144 (98%)	-0.36	0 100 100	13, 23, 44, 53	0
1	f	142/144 (98%)	-0.30	1 (0%) 87 87	13, 24, 42, 49	0
1	g	142/144 (98%)	-0.31	1 (0%) 87 87	14, 23, 42, 57	1 (0%)
1	h	142/144 (98%)	-0.38	0 100 100	11, 21, 40, 48	1 (0%)
1	i	142/144 (98%)	-0.40	2 (1%) 75 75	13, 21, 39, 46	1 (0%)
1	j	143/144 (99%)	-0.37	0 100 100	13, 22, 43, 51	1 (0%)
1	k	142/144 (98%)	-0.31	1 (0%) 87 87	15, 27, 45, 58	1 (0%)
1	l	142/144 (98%)	-0.20	1 (0%) 87 87	16, 28, 50, 63	0
1	m	142/144 (98%)	-0.27	0 100 100	16, 28, 49, 57	0
All	All	6822/6912 (98%)	-0.34	18 (0%) 93 93	11, 24, 43, 66	49 (0%)

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	34	VAL	3.0
1	A	99	LEU	2.9
1	C	101	ASP	2.6
1	g	101	ASP	2.5
1	5	99	LEU	2.5

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Mol	Chain	Res	Type	RSRZ
1	Z	40	GLU	2.5
1	F	99	LEU	2.2
1	l	13	THR	2.2
1	1	99	LEU	2.2
1	i	133	GLU	2.2
1	H	99	LEU	2.2
1	k	79	GLY	2.1
1	O	131	GLY	2.1
1	i	131	GLY	2.0
1	Z	99	LEU	2.0
1	d	129	HIS	2.0
1	f	12	GLU	2.0
1	M	61	VAL	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
5	PO4	l	1148	5/5	0.89	0.37	12.68	40,46,51,61	0
4	MG	J	1147	1/1	0.99	0.11	0.76	17,17,17,17	0
4	MG	g	1147	1/1	0.99	0.11	0.52	16,16,16,16	0
3	POP	V	1146	9/9	0.99	0.09	0.48	14,16,17,20	0
5	PO4	e	1148	5/5	0.94	0.14	0.42	49,54,55,61	0
2	DUR	3	1145	16/16	0.97	0.12	0.39	12,13,15,16	0
4	MG	P	1147	1/1	0.99	0.11	0.38	13,13,13,13	0
4	MG	U	1147	1/1	0.99	0.09	0.38	15,15,15,15	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
4	MG	m	1147	1/1	0.99	0.10	0.25	17,17,17,17	0
2	DUR	J	1145	16/16	0.96	0.10	0.19	13,14,16,16	0
2	DUR	L	1145	16/16	0.98	0.10	0.19	14,15,18,19	0
2	DUR	b	1145	16/16	0.98	0.11	0.19	13,15,17,19	0
2	DUR	N	1145	16/16	0.98	0.12	0.14	15,17,20,20	0
2	DUR	W	1145	16/16	0.97	0.11	0.13	14,15,17,17	0
2	DUR	9	1145	16/16	0.98	0.10	0.13	13,14,16,17	0
4	MG	R	1147	1/1	0.99	0.10	0.11	12,12,12,12	0
4	MG	7	1147	1/1	0.98	0.10	0.09	20,20,20,20	0
4	MG	b	1147	1/1	0.99	0.09	0.09	16,16,16,16	0
2	DUR	M	1145	16/16	0.98	0.11	0.07	14,17,18,19	0
2	DUR	Y	1145	16/16	0.95	0.12	0.07	14,17,20,20	0
2	DUR	O	1145	16/16	0.98	0.10	0.06	16,18,21,21	0
2	DUR	C	1145	16/16	0.96	0.11	0.03	16,18,22,22	0
3	POP	Z	1146	9/9	0.99	0.10	0.02	16,18,19,20	0
2	DUR	K	1145	16/16	0.97	0.11	0.02	14,16,18,18	0
2	DUR	d	1145	16/16	0.98	0.10	-0.02	14,15,17,18	0
3	POP	a	1146	9/9	0.99	0.09	-0.08	13,15,16,17	0
3	POP	7	1146	9/9	0.99	0.09	-0.10	18,20,22,23	0
2	DUR	A	1145	16/16	0.98	0.10	-0.12	16,18,20,20	0
2	DUR	Z	1145	16/16	0.97	0.10	-0.17	17,19,22,22	0
2	DUR	j	1145	16/16	0.97	0.11	-0.18	14,16,17,20	0
2	DUR	Q	1145	16/16	0.97	0.10	-0.19	14,15,17,20	0
4	MG	D	1147	1/1	0.99	0.08	-0.21	15,15,15,15	0
3	POP	m	1146	9/9	0.99	0.08	-0.23	18,20,21,22	0
2	DUR	i	1145	16/16	0.97	0.09	-0.24	15,16,19,19	0
2	DUR	l	1145	16/16	0.97	0.10	-0.25	17,20,23,23	0
2	DUR	G	1145	16/16	0.98	0.10	-0.26	11,13,14,14	0
2	DUR	I	1145	16/16	0.97	0.09	-0.28	10,13,14,15	0
2	DUR	c	1145	16/16	0.97	0.09	-0.28	16,17,19,21	0
2	DUR	B	1145	16/16	0.97	0.10	-0.30	18,19,20,20	0
2	DUR	T	1145	16/16	0.97	0.09	-0.30	14,18,20,20	0
4	MG	9	1147	1/1	0.99	0.08	-0.31	17,17,17,17	0
3	POP	2	1146	9/9	0.99	0.08	-0.31	13,14,16,17	0
3	POP	N	1146	9/9	0.99	0.08	-0.33	16,18,19,20	0
2	DUR	6	1145	16/16	0.98	0.10	-0.33	19,21,23,23	0
3	POP	g	1146	9/9	0.99	0.08	-0.36	13,15,16,19	0
3	POP	I	1146	9/9	0.99	0.08	-0.36	11,13,15,17	0
2	DUR	E	1145	16/16	0.98	0.10	-0.37	13,14,15,15	0
2	DUR	8	1145	16/16	0.97	0.09	-0.41	12,13,14,15	0
4	MG	A	1147	1/1	0.99	0.09	-0.42	21,21,21,21	0
2	DUR	h	1145	16/16	0.97	0.09	-0.43	11,13,14,14	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
4	MG	3	1147	1/1	0.98	0.08	-0.43	17,17,17,17	0
3	POP	i	1146	9/9	1.00	0.09	-0.43	13,16,19,21	0
2	DUR	k	1145	16/16	0.97	0.10	-0.44	17,18,20,20	0
2	DUR	e	1145	16/16	0.97	0.09	-0.45	14,15,16,16	0
2	DUR	S	1145	16/16	0.97	0.09	-0.46	12,15,17,17	0
2	DUR	D	1145	16/16	0.97	0.09	-0.46	13,14,17,20	0
3	POP	L	1146	9/9	0.99	0.09	-0.48	16,17,19,21	0
2	DUR	P	1145	16/16	0.98	0.09	-0.48	13,14,16,17	0
4	MG	H	1147	1/1	0.97	0.08	-0.55	16,16,16,16	0
3	POP	3	1146	9/9	0.99	0.08	-0.55	14,16,17,18	0
4	MG	S	1147	1/1	0.99	0.08	-0.57	14,14,14,14	0
3	POP	W	1146	9/9	0.99	0.08	-0.58	10,13,15,16	0
3	POP	j	1146	9/9	0.99	0.08	-0.58	13,14,16,17	0
3	POP	Q	1146	9/9	0.99	0.08	-0.59	11,12,14,14	0
2	DUR	2	1145	16/16	0.97	0.08	-0.60	12,14,17,18	0
2	DUR	g	1145	16/16	0.97	0.09	-0.62	14,16,18,19	0
2	DUR	U	1145	16/16	0.97	0.09	-0.62	15,16,17,18	0
2	DUR	V	1145	16/16	0.97	0.09	-0.63	12,14,16,17	0
3	POP	C	1146	9/9	0.99	0.07	-0.64	17,18,22,22	0
2	DUR	1	1145	16/16	0.95	0.08	-0.66	14,18,19,20	0
3	POP	R	1146	9/9	0.99	0.08	-0.66	13,14,15,16	0
2	DUR	f	1145	16/16	0.98	0.09	-0.67	15,16,19,22	0
2	DUR	4	1145	16/16	0.97	0.08	-0.68	11,12,13,14	0
3	POP	f	1146	9/9	1.00	0.07	-0.70	13,15,17,17	0
4	MG	8	1147	1/1	0.99	0.07	-0.74	17,17,17,17	0
3	POP	F	1146	9/9	0.99	0.07	-0.75	11,13,13,13	0
3	POP	U	1146	9/9	0.99	0.08	-0.75	16,17,19,19	0
3	POP	D	1146	9/9	0.99	0.07	-0.77	11,15,17,17	0
4	MG	k	1147	1/1	0.98	0.09	-0.81	21,21,21,21	0
3	POP	4	1146	9/9	0.99	0.08	-0.84	14,14,16,17	0
2	DUR	R	1145	16/16	0.97	0.08	-0.85	12,13,15,15	0
3	POP	T	1146	9/9	0.99	0.07	-0.86	13,16,18,19	0
2	DUR	H	1145	16/16	0.98	0.07	-0.88	12,14,15,15	0
2	DUR	m	1145	16/16	0.98	0.08	-0.90	15,20,23,23	0
3	POP	h	1146	9/9	0.99	0.07	-0.91	17,18,19,19	0
3	POP	K	1146	9/9	0.99	0.07	-0.94	13,15,18,21	0
4	MG	X	1147	1/1	0.99	0.08	-0.95	16,16,16,16	0
2	DUR	F	1145	16/16	0.97	0.07	-0.95	12,14,15,16	0
3	POP	P	1146	9/9	0.99	0.07	-0.98	14,15,17,17	0
3	POP	X	1146	9/9	0.99	0.07	-1.00	15,16,18,19	0
4	MG	V	1147	1/1	0.99	0.07	-1.02	14,14,14,14	0
2	DUR	a	1145	16/16	0.98	0.07	-1.02	15,16,17,17	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
4	MG	6	1147	1/1	0.98	0.08	-1.02	19,19,19,19	0
3	POP	e	1146	9/9	0.99	0.08	-1.03	16,18,20,20	0
2	DUR	X	1145	16/16	0.98	0.07	-1.11	12,14,15,20	0
2	DUR	5	1145	16/16	0.98	0.07	-1.13	17,20,22,22	0
4	MG	C	1147	1/1	0.98	0.06	-1.13	13,13,13,13	0
3	POP	B	1146	9/9	0.99	0.06	-1.13	16,18,20,23	0
4	MG	j	1147	1/1	0.98	0.07	-1.14	13,13,13,13	0
3	POP	E	1146	9/9	0.99	0.06	-1.15	15,16,17,18	0
4	MG	5	1147	1/1	0.99	0.08	-1.18	18,18,18,18	0
3	POP	Y	1146	9/9	0.99	0.06	-1.19	15,17,18,19	0
3	POP	9	1146	9/9	0.99	0.07	-1.19	14,14,16,16	0
3	POP	J	1146	9/9	0.99	0.07	-1.19	15,18,19,20	0
3	POP	M	1146	9/9	0.99	0.06	-1.23	13,19,20,20	0
3	POP	l	1146	9/9	0.99	0.06	-1.24	15,19,22,23	0
3	POP	G	1146	9/9	0.99	0.06	-1.26	12,14,15,16	0
3	POP	O	1146	9/9	0.99	0.07	-1.28	16,17,22,23	0
2	DUR	7	1145	16/16	0.98	0.08	-1.28	15,18,21,24	0
3	POP	A	1146	9/9	0.99	0.07	-1.36	18,19,21,25	0
3	POP	6	1146	9/9	0.99	0.07	-1.36	18,21,22,24	0
4	MG	4	1147	1/1	0.97	0.06	-1.37	15,15,15,15	0
3	POP	5	1146	9/9	0.99	0.07	-1.38	17,19,20,20	0
3	POP	1	1146	9/9	0.99	0.06	-1.40	18,19,21,24	0
3	POP	8	1146	9/9	0.99	0.06	-1.41	14,16,17,17	0
3	POP	k	1146	9/9	0.99	0.07	-1.41	19,23,24,25	0
4	MG	Q	1147	1/1	0.99	0.06	-1.49	10,10,10,10	0
4	MG	K	1147	1/1	0.99	0.07	-1.53	15,15,15,15	0
3	POP	H	1146	9/9	0.99	0.07	-1.55	14,15,16,17	0
4	MG	F	1147	1/1	0.98	0.06	-1.55	14,14,14,14	0
4	MG	E	1147	1/1	0.99	0.06	-1.59	17,17,17,17	0
3	POP	d	1146	9/9	0.99	0.07	-1.60	19,20,22,23	0
3	POP	c	1146	9/9	0.99	0.06	-1.63	14,16,18,18	0
4	MG	N	1147	1/1	0.99	0.05	-1.73	14,14,14,14	0
3	POP	b	1146	9/9	0.99	0.06	-1.73	15,17,20,21	0
4	MG	c	1147	1/1	0.99	0.06	-1.76	15,15,15,15	0
4	MG	1	1147	1/1	0.99	0.05	-1.83	18,18,18,18	0
4	MG	L	1147	1/1	1.00	0.05	-1.83	17,17,17,17	0
3	POP	S	1146	9/9	0.99	0.06	-1.85	13,15,18,21	0
4	MG	Z	1147	1/1	0.97	0.05	-1.86	21,21,21,21	0
4	MG	M	1147	1/1	0.95	0.04	-2.03	19,19,19,19	0
4	MG	T	1147	1/1	0.99	0.06	-2.10	17,17,17,17	0
4	MG	I	1147	1/1	0.99	0.04	-2.17	11,11,11,11	0
4	MG	l	1147	1/1	0.99	0.04	-2.22	18,18,18,18	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
4	MG	2	1147	1/1	0.99	0.04	-2.26	11,11,11,11	0
4	MG	i	1147	1/1	1.00	0.05	-2.31	14,14,14,14	0
4	MG	W	1147	1/1	0.99	0.06	-2.46	12,12,12,12	0
4	MG	O	1147	1/1	0.99	0.04	-2.53	20,20,20,20	0
4	MG	G	1147	1/1	0.99	0.04	-2.58	14,14,14,14	0
4	MG	f	1147	1/1	0.99	0.05	-2.81	17,17,17,17	0
4	MG	h	1147	1/1	0.99	0.03	-2.86	15,15,15,15	0
4	MG	B	1147	1/1	0.97	0.04	-3.00	18,18,18,18	0
4	MG	a	1147	1/1	0.99	0.05	-3.19	15,15,15,15	0
4	MG	e	1147	1/1	0.99	0.04	-3.37	18,18,18,18	0
4	MG	d	1147	1/1	0.96	0.06	-3.48	22,22,22,22	0
4	MG	Y	1147	1/1	0.99	0.03	-3.57	16,16,16,16	0

6.5 Other polymers [i](#)

There are no such residues in this entry.