



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 02:16 AM UTC

PDB ID : 3OL8 / pdb_00003ol8
Title : Poliovirus polymerase elongation complex with CTP-Mn
Authors : Gong, P.; Peersen, O.B.
Deposited on : 2010-08-25
Resolution : 2.75 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

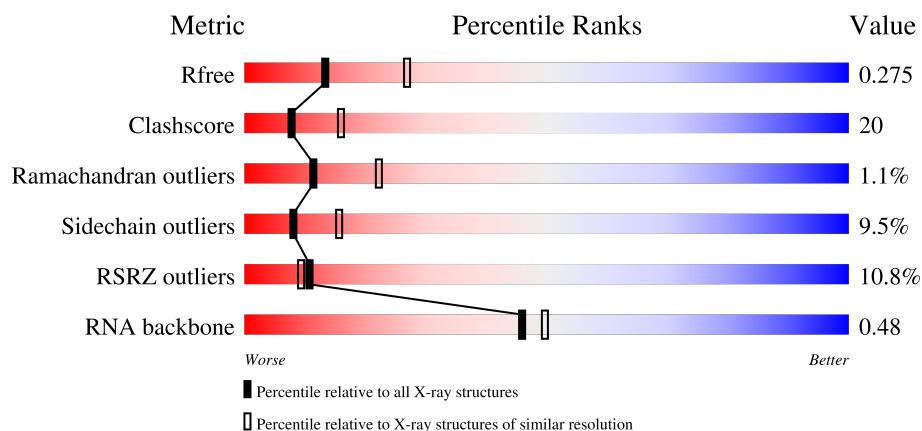
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.75 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1009 (2.76-2.76)
Clashscore	190562	1044 (2.76-2.76)
Ramachandran outliers	187476	1024 (2.76-2.76)
Sidechain outliers	187428	1024 (2.76-2.76)
RSRZ outliers	180081	1009 (2.76-2.76)
RNA backbone	3983	1179 (3.00-2.52)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	471	8% 62% 30% 6%
1	E	471	8% 62% 30% 6%
1	I	471	11% 56% 37% 5%
1	M	471	13% 55% 38% 5%

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Mol	Chain	Length	Quality of chain
2	B	26	
2	F	26	
2	J	26	
2	N	26	
3	C	15	
3	G	15	
3	K	15	
3	O	15	
4	D	9	
4	H	9	
4	L	9	
4	P	9	

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
8	IPA	A	6011	-	-	X	-
8	IPA	E	6014	-	-	X	-
8	IPA	M	6019	-	-	X	-

2 Entry composition

There are 10 unique types of molecules in this entry. The entry contains 18318 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 Polymerase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	461	Total	C	N	O	S	0	0	0
			3697	2370	610	695	22			
1	E	461	Total	C	N	O	S	0	0	0
			3697	2370	610	695	22			
1	I	461	Total	C	N	O	S	0	0	0
			3697	2370	610	695	22			
1	M	461	Total	C	N	O	S	0	0	0
			3697	2370	610	695	22			

There are 44 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	446	ASP	LEU	engineered mutation	UNP B3VQP5
A	462	GLY	-	expression tag	UNP B3VQP5
A	463	SER	-	expression tag	UNP B3VQP5
A	464	SER	-	expression tag	UNP B3VQP5
A	465	SER	-	expression tag	UNP B3VQP5
A	466	HIS	-	expression tag	UNP B3VQP5
A	467	HIS	-	expression tag	UNP B3VQP5
A	468	HIS	-	expression tag	UNP B3VQP5
A	469	HIS	-	expression tag	UNP B3VQP5
A	470	HIS	-	expression tag	UNP B3VQP5
A	471	HIS	-	expression tag	UNP B3VQP5
E	446	ASP	LEU	engineered mutation	UNP B3VQP5
E	462	GLY	-	expression tag	UNP B3VQP5
E	463	SER	-	expression tag	UNP B3VQP5
E	464	SER	-	expression tag	UNP B3VQP5
E	465	SER	-	expression tag	UNP B3VQP5
E	466	HIS	-	expression tag	UNP B3VQP5
E	467	HIS	-	expression tag	UNP B3VQP5
E	468	HIS	-	expression tag	UNP B3VQP5
E	469	HIS	-	expression tag	UNP B3VQP5
E	470	HIS	-	expression tag	UNP B3VQP5

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Chain	Residue	Modelled	Actual	Comment	Reference
E	471	HIS	-	expression tag	UNP B3VQP5
I	446	ASP	LEU	engineered mutation	UNP B3VQP5
I	462	GLY	-	expression tag	UNP B3VQP5
I	463	SER	-	expression tag	UNP B3VQP5
I	464	SER	-	expression tag	UNP B3VQP5
I	465	SER	-	expression tag	UNP B3VQP5
I	466	HIS	-	expression tag	UNP B3VQP5
I	467	HIS	-	expression tag	UNP B3VQP5
I	468	HIS	-	expression tag	UNP B3VQP5
I	469	HIS	-	expression tag	UNP B3VQP5
I	470	HIS	-	expression tag	UNP B3VQP5
I	471	HIS	-	expression tag	UNP B3VQP5
M	446	ASP	LEU	engineered mutation	UNP B3VQP5
M	462	GLY	-	expression tag	UNP B3VQP5
M	463	SER	-	expression tag	UNP B3VQP5
M	464	SER	-	expression tag	UNP B3VQP5
M	465	SER	-	expression tag	UNP B3VQP5
M	466	HIS	-	expression tag	UNP B3VQP5
M	467	HIS	-	expression tag	UNP B3VQP5
M	468	HIS	-	expression tag	UNP B3VQP5
M	469	HIS	-	expression tag	UNP B3VQP5
M	470	HIS	-	expression tag	UNP B3VQP5
M	471	HIS	-	expression tag	UNP B3VQP5

- Molecule 2 is a RNA chain called RNA (5'-R(*AP*AP*GP*UP*CP*UP*CP*CP*AP*GP*GP*UP*CP*UP*CP*UP*CP*GP*UP*CP*CP*GP*GP*AP*AP*A)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	17	Total	C	N	O	P	0	0	0
			341	150	56	118	17			
2	F	17	Total	C	N	O	P	0	0	0
			341	150	56	118	17			
2	J	17	Total	C	N	O	P	0	0	0
			341	150	56	118	17			
2	N	17	Total	C	N	O	P	0	0	0
			341	150	56	118	17			

- Molecule 3 is a RNA chain called RNA (5'-R(*GP*CP*CP*CP*GP*GP*AP*CP*GP*AP*GP*AP*GP*AP*C)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	15	Total	C	N	O	P	0	0	0
			323	145	65	99	14			
3	G	15	Total	C	N	O	P	0	0	0
			323	145	65	99	14			
3	K	15	Total	C	N	O	P	0	0	0
			323	145	65	99	14			
3	O	15	Total	C	N	O	P	0	0	0
			323	145	65	99	14			

- Molecule 4 is a RNA chain called RNA (5'-R(*GP*GP*GP*AP*GP*AP*UP*GP*A)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	3	Total	C	N	O	P	0	0	0
			68	30	15	20	3			
4	H	3	Total	C	N	O	P	0	0	0
			68	30	15	20	3			
4	L	4	Total	C	N	O	P	0	0	0
			91	40	20	27	4			
4	P	4	Total	C	N	O	P	0	0	0
			91	40	20	27	4			

- Molecule 5 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	Zn	0	0
			1	1		
5	E	1	Total	Zn	0	0
			1	1		
5	I	1	Total	Zn	0	0
			1	1		
5	M	1	Total	Zn	0	0
			1	1		

- Molecule 6 is MANGANESE (II) ION (CCD ID: MN) (formula: Mn).

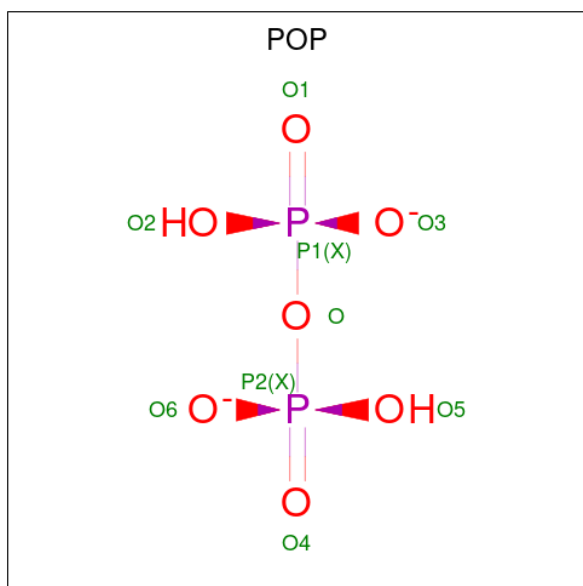
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	2	Total	Mn	0	0
			2	2		
6	E	2	Total	Mn	0	0
			2	2		
6	I	2	Total	Mn	0	0
			2	2		

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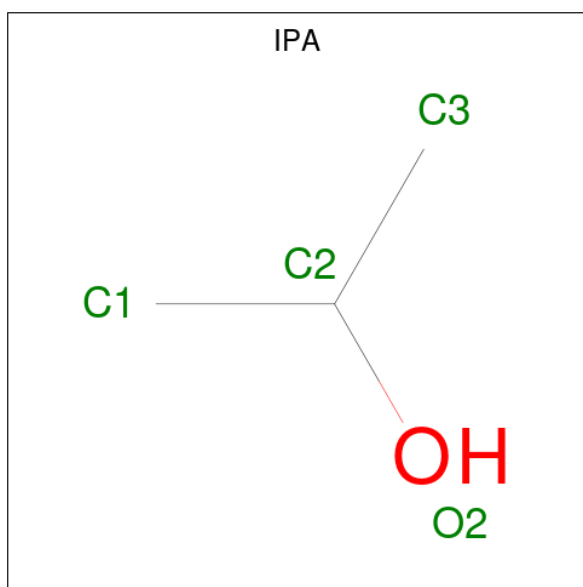
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	M	2	Total	Mn	0	0
			2	2		

- Molecule 7 is PYROPHOSPHATE 2- (CCD ID: POP) (formula: $\text{H}_2\text{O}_7\text{P}_2$).



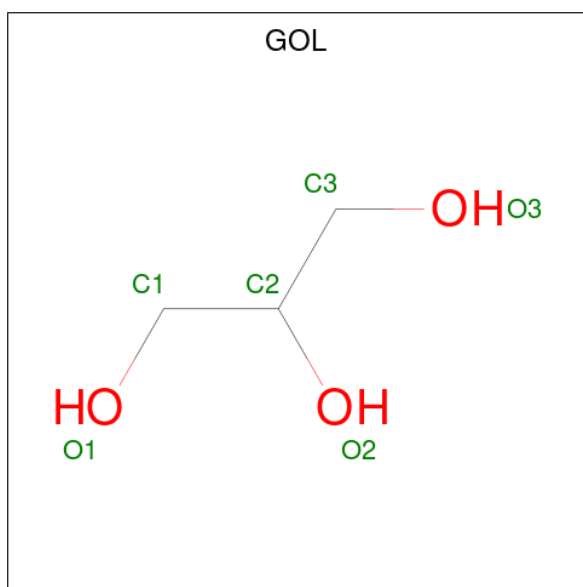
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	O	P	0	0
			9	7	2		
7	G	1	Total	O	P	0	0
			9	7	2		
7	I	1	Total	O	P	0	0
			9	7	2		
7	O	1	Total	O	P	0	0
			9	7	2		

- Molecule 8 is ISOPROPYL ALCOHOL (CCD ID: IPA) (formula: $\text{C}_3\text{H}_8\text{O}$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	A	1	Total	C	O	0	0
			4	3	1		
8	A	1	Total	C	O	0	0
			4	3	1		
8	A	1	Total	C	O	0	0
			4	3	1		
8	A	1	Total	C	O	0	0
			4	3	1		
8	A	1	Total	C	O	0	0
			4	3	1		
8	E	1	Total	C	O	0	0
			4	3	1		
8	E	1	Total	C	O	0	0
			4	3	1		
8	M	1	Total	C	O	0	0
			4	3	1		
8	M	1	Total	C	O	0	0
			4	3	1		

- Molecule 9 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
9	A	1	Total	C	O	0	0
			6	3	3		
9	E	1	Total	C	O	0	0
			6	3	3		
9	J	1	Total	C	O	0	0
			6	3	3		
9	N	1	Total	C	O	0	0
			6	3	3		

- Molecule 10 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	A	106	Total	O	0	0
			106	106		
10	B	12	Total	O	0	0
			12	12		
10	C	7	Total	O	0	0
			7	7		
10	E	111	Total	O	0	0
			111	111		
10	F	13	Total	O	0	0
			13	13		
10	G	8	Total	O	0	0
			8	8		
10	I	76	Total	O	0	0
			76	76		
10	J	14	Total	O	0	0
			14	14		

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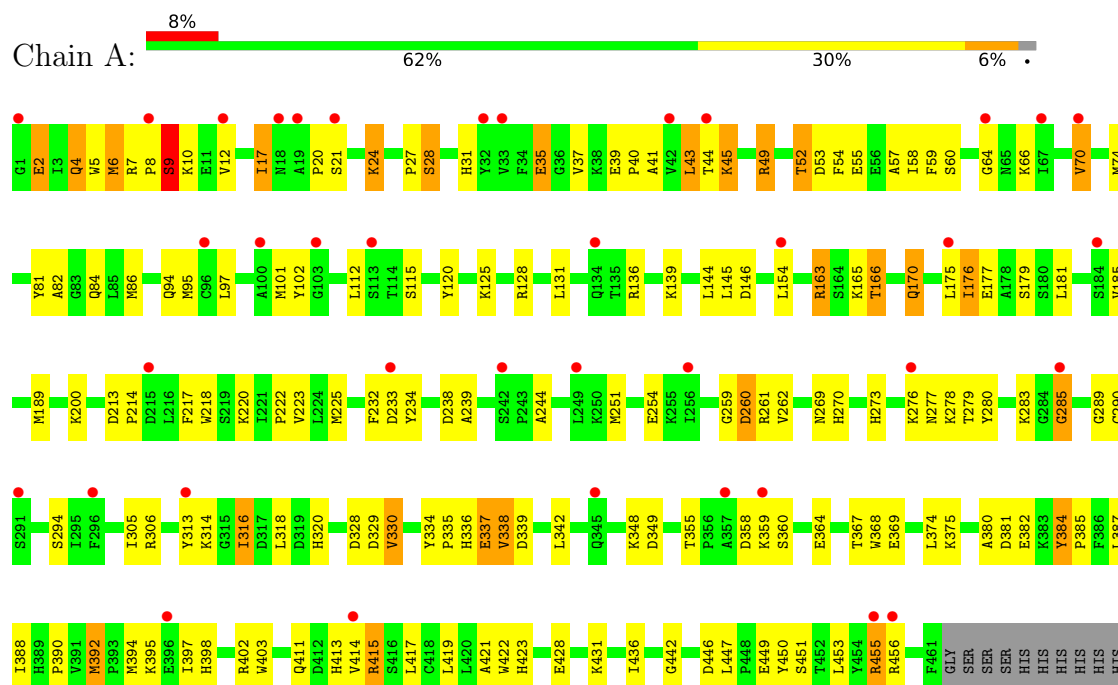
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	K	5	Total 5	O 5	0	0
10	M	74	Total 74	O 74	0	0
10	N	16	Total 16	O 16	0	0
10	O	4	Total 4	O 4	0	0
10	P	2	Total 2	O 2	0	0

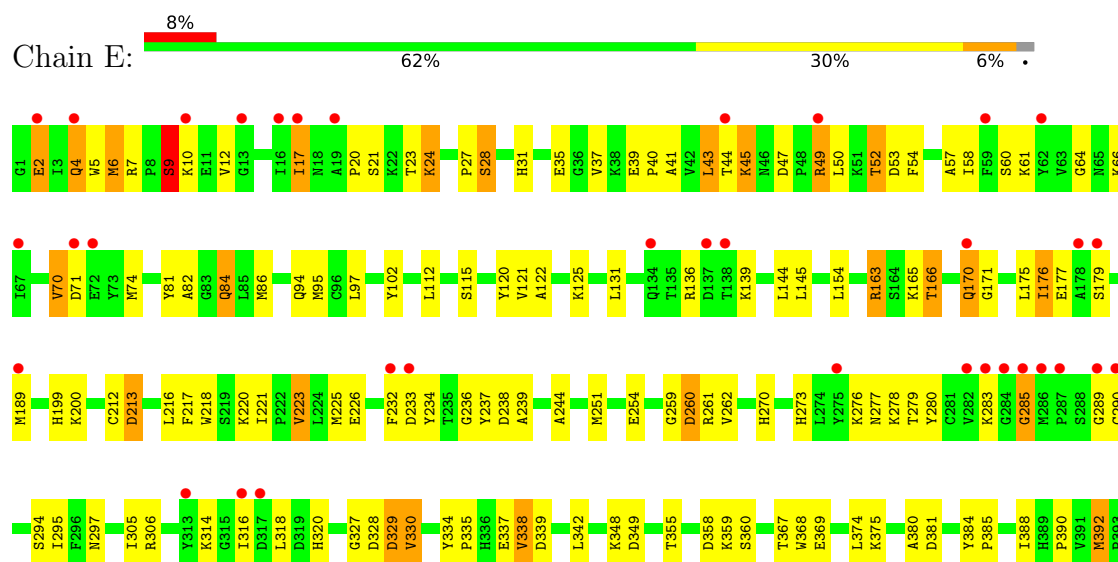
3 Residue-property plots

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

• Molecule 1: Polymerase

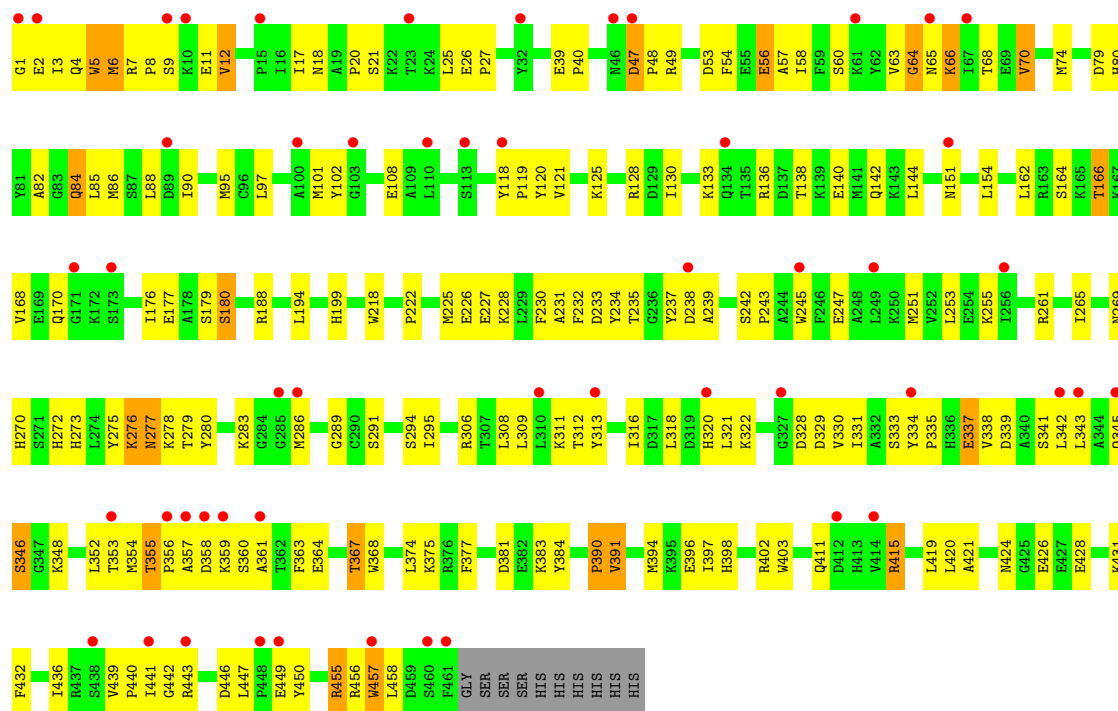


• Molecule 1: Polymerase

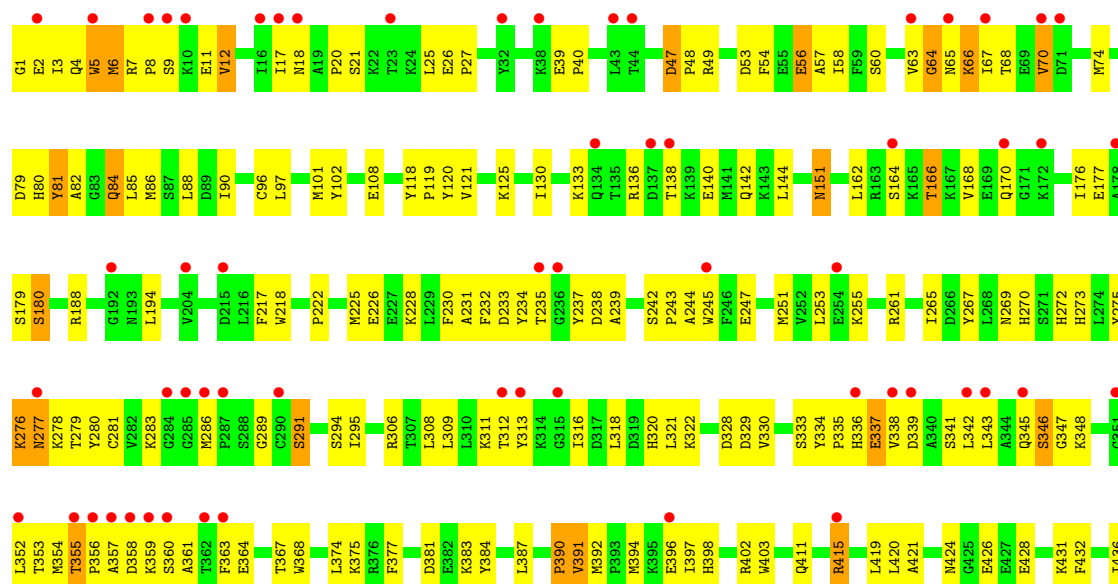




• Molecule 1: Polymerase

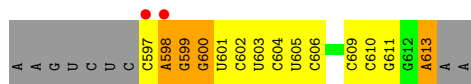
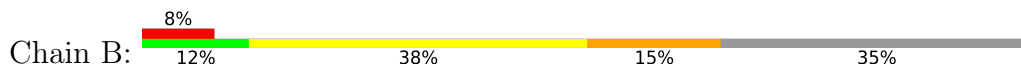


• Molecule 1: Polymerase

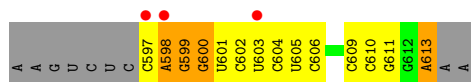




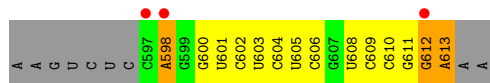
- Molecule 2: RNA (5'-R(*AP*AP*GP*UP*CP*UP*CP*CP*AP*GP*GP*UP*CP*UP*CP*UP*CP*GP*UP*CP*CP*GP*GP*AP*AP*A)-3')



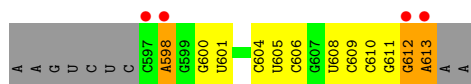
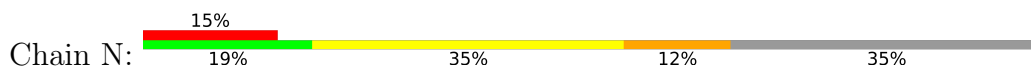
- Molecule 2: RNA (5'-R(*AP*AP*GP*UP*CP*UP*CP*CP*AP*GP*GP*UP*CP*UP*CP*UP*CP*GP*UP*CP*CP*GP*GP*AP*AP*A)-3')



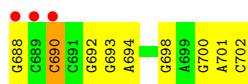
- Molecule 2: RNA (5'-R(*AP*AP*GP*UP*CP*UP*CP*CP*AP*GP*GP*UP*CP*UP*CP*UP*CP*GP*UP*CP*CP*GP*GP*AP*AP*A)-3')



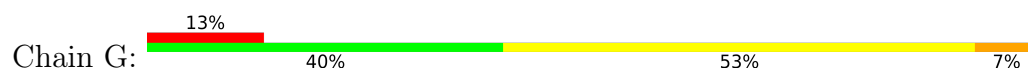
- Molecule 2: RNA (5'-R(*AP*AP*GP*UP*CP*UP*CP*CP*AP*GP*GP*UP*CP*UP*CP*UP*CP*GP*UP*CP*CP*GP*GP*AP*AP*A)-3')



- Molecule 3: RNA (5'-R(*GP*CP*CP*CP*GP*GP*AP*CP*GP*AP*GP*AP*GP*AP*C)-3')



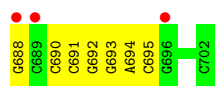
- Molecule 3: RNA (5'-R(*GP*CP*CP*CP*GP*GP*AP*CP*GP*AP*GP*AP*GP*AP*C)-3')



- Molecule 3: RNA (5'-R(*GP*CP*CP*CP*GP*GP*AP*CP*GP*AP*GP*AP*GP*AP*C)-3')



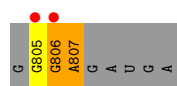
- Molecule 3: RNA (5'-R(*GP*CP*CP*CP*GP*GP*AP*CP*GP*AP*GP*AP*GP*AP*C)-3')



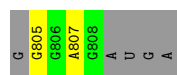
- Molecule 4: RNA (5'-R(*GP*GP*GP*AP*GP*AP*UP*GP*A)-3')



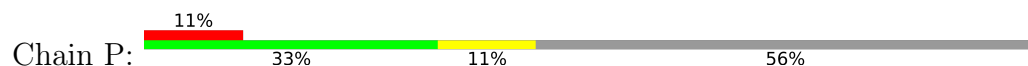
- Molecule 4: RNA (5'-R(*GP*GP*GP*AP*GP*AP*UP*GP*A)-3')



- Molecule 4: RNA (5'-R(*GP*GP*GP*AP*GP*AP*UP*GP*A)-3')



- Molecule 4: RNA (5'-R(*GP*GP*GP*AP*GP*AP*UP*GP*A)-3')



G	G805	G806	A807	G808	A	U	G	A
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4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	60.78Å 60.86Å 193.08Å 83.06° 83.04° 76.81°	Depositor
Resolution (Å)	44.16 – 2.75 44.16 – 2.75	Depositor EDS
% Data completeness (in resolution range)	97.0 (44.16-2.75) 97.5 (44.16-2.75)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.16 (at 2.39Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.6.4_486)	Depositor
R, R_{free}	0.224 , 0.277 0.219 , 0.275	Depositor DCC
R_{free} test set	5134 reflections (4.93%)	wwPDB-VP
Wilson B-factor (Å ²)	67.0	Xtriage
Anisotropy	0.337	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 92.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.63$, $\langle L^2 \rangle = 0.49$	Xtriage
Estimated twinning fraction	0.398 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	18318	wwPDB-VP
Average B, all atoms (Å ²)	72.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 45.86 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.2443e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, IPA, POP, GOL, MN

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# $ Z > 5$	RMSZ	# $ Z > 5$
1	A	0.57	0/3787	0.91	6/5122 (0.1%)
1	E	0.58	0/3787	0.90	5/5122 (0.1%)
1	I	0.56	0/3787	0.87	7/5122 (0.1%)
1	M	0.57	0/3787	0.87	7/5122 (0.1%)
2	B	0.59	2/378 (0.5%)	0.57	0/587
2	F	0.58	2/378 (0.5%)	0.56	0/587
2	J	0.59	2/378 (0.5%)	0.49	0/587
2	N	0.59	2/378 (0.5%)	0.49	0/587
3	C	0.38	0/362	0.44	0/564
3	G	0.37	0/362	0.44	0/564
3	K	0.36	0/362	0.44	0/564
3	O	0.36	0/362	0.43	0/564
4	D	0.22	0/76	0.30	0/117
4	H	0.19	0/76	0.29	0/117
4	L	0.25	0/102	0.30	0/158
4	P	0.23	0/102	0.30	0/158
All	All	0.55	8/18464 (0.0%)	0.82	25/25642 (0.1%)

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	N	613	A	P-OP1	6.68	1.61	1.48
2	J	613	A	P-OP1	6.30	1.61	1.48
2	B	613	A	P-OP1	6.15	1.60	1.48
2	F	613	A	P-OP1	6.00	1.60	1.48
2	B	613	A	P-OP2	5.83	1.60	1.48
2	F	613	A	P-OP2	5.71	1.59	1.48
2	J	613	A	P-OP2	5.56	1.59	1.48
2	N	613	A	P-OP2	5.17	1.58	1.48

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	I	390	PRO	CB-CA-C	-6.11	106.74	111.87
1	A	213	ASP	CA-C-N	5.96	126.12	119.32
1	A	213	ASP	C-N-CA	5.96	126.12	119.32
1	M	47	ASP	CA-C-N	5.95	125.42	119.24
1	M	47	ASP	C-N-CA	5.95	125.42	119.24
1	M	233	ASP	N-CA-C	-5.77	100.63	109.41
1	M	390	PRO	CB-CA-C	-5.75	107.04	111.87
1	I	47	ASP	CA-C-N	5.70	125.17	119.24
1	I	47	ASP	C-N-CA	5.70	125.17	119.24
1	I	5	TRP	N-CA-C	5.66	115.75	108.34
1	M	5	TRP	N-CA-C	5.53	115.58	108.34
1	I	130	ILE	CB-CA-C	-5.49	106.03	111.30
1	E	213	ASP	CA-C-N	5.44	125.53	119.32
1	E	213	ASP	C-N-CA	5.44	125.53	119.32
1	M	130	ILE	CB-CA-C	-5.36	106.16	111.30
1	A	384	TYR	CA-C-N	5.31	125.00	119.64
1	A	384	TYR	C-N-CA	5.31	125.00	119.64
1	E	9	SER	N-CA-C	-5.28	106.81	113.20
1	E	233	ASP	N-CA-C	-5.16	102.26	109.69
1	A	9	SER	N-CA-C	-5.13	106.99	113.20
1	M	81	TYR	N-CA-C	-5.07	105.65	111.07
1	A	233	ASP	N-CA-C	-5.03	102.45	109.69
1	E	221	ILE	CB-CA-C	-5.02	109.03	114.35
1	I	233	ASP	N-CA-C	-5.02	101.78	109.41
1	I	227	GLU	N-CA-C	5.01	117.39	111.33

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3697	0	3658	128	0
1	E	3697	0	3658	127	0
1	I	3697	0	3658	159	0
1	M	3697	0	3658	178	0
2	B	341	0	172	23	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	F	341	0	172	17	0
2	J	341	0	172	15	0
2	N	341	0	172	13	0
3	C	323	0	167	12	0
3	G	323	0	167	14	0
3	K	323	0	167	3	0
3	O	323	0	167	5	0
4	D	68	0	34	6	0
4	H	68	0	34	4	0
4	L	91	0	45	1	0
4	P	91	0	45	0	0
5	A	1	0	0	0	0
5	E	1	0	0	0	0
5	I	1	0	0	0	0
5	M	1	0	0	0	0
6	A	2	0	0	0	0
6	E	2	0	0	0	0
6	I	2	0	0	0	0
6	M	2	0	0	0	0
7	A	9	0	0	0	0
7	G	9	0	0	1	0
7	I	9	0	0	1	0
7	O	9	0	0	0	0
8	A	20	0	40	12	0
8	E	8	0	16	14	0
8	M	8	0	16	6	0
9	A	6	0	8	0	0
9	E	6	0	8	0	0
9	J	6	0	8	0	0
9	N	6	0	8	0	0
10	A	106	0	0	7	0
10	B	12	0	0	1	0
10	C	7	0	0	2	0
10	E	111	0	0	9	0
10	F	13	0	0	0	0
10	G	8	0	0	1	0
10	I	76	0	0	6	0
10	J	14	0	0	0	0
10	K	5	0	0	0	0
10	M	74	0	0	5	0
10	N	16	0	0	1	0
10	O	4	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
10	P	2	0	0	0	0
All	All	18318	0	16250	678	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 20.

All (678) 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:M:232:PHE:CE1	1:M:354:MET:HE2	1.97	0.99
1:I:232:PHE:CE1	1:I:354:MET:HE2	1.98	0.99
1:M:247:GLU:O	1:M:251:MET:HG3	1.64	0.97
1:I:247:GLU:O	1:I:251:MET:HG3	1.65	0.96
1:E:368:TRP:HB3	8:E:6029:IPA:H2	1.49	0.94
1:A:128:ARG:HG3	10:A:564:HOH:O	1.66	0.93
2:B:613:A:O5'	3:G:688:G:H5'	1.68	0.93
1:E:217:PHE:HB2	8:E:6014:IPA:H32	1.53	0.89
1:I:311:LYS:HD3	1:I:346:SER:HB3	1.55	0.88
1:M:311:LYS:HD3	1:M:346:SER:HB3	1.55	0.87
1:A:70:VAL:HG11	1:A:251:MET:HE1	1.57	0.85
1:M:309:LEU:HD23	1:M:343:LEU:HD21	1.57	0.85
3:C:688:G:H5'	2:F:613:A:O5'	1.75	0.84
1:I:309:LEU:HD23	1:I:343:LEU:HD21	1.57	0.84
1:M:243:PRO:HB3	10:M:492:HOH:O	1.77	0.83
1:I:270:HIS:HD2	1:I:283:LYS:HG2	1.42	0.83
1:E:6:MET:HG2	1:E:280:TYR:HB3	1.60	0.82
1:M:377:PHE:HB2	1:M:391:VAL:HG22	1.61	0.82
2:B:598:A:H1'	4:D:805:G:N2	1.94	0.82
1:E:9:SER:HB3	1:E:277:ASN:O	1.79	0.82
1:E:254:GLU:HG3	1:E:262:VAL:HG11	1.61	0.82
1:A:9:SER:HB3	1:A:277:ASN:O	1.78	0.82
1:A:6:MET:HG2	1:A:280:TYR:HB3	1.60	0.82
1:M:7:ARG:HD2	1:M:12:VAL:HG23	1.63	0.80
1:M:270:HIS:HD2	1:M:283:LYS:HG2	1.44	0.80
1:A:254:GLU:HG3	1:A:262:VAL:HG11	1.63	0.80
1:I:7:ARG:HD2	1:I:12:VAL:HG23	1.64	0.79
1:E:70:VAL:HG11	1:E:251:MET:HE1	1.62	0.79
1:A:9:SER:HB2	1:A:279:THR:OG1	1.83	0.79
1:A:270:HIS:HD2	1:A:283:LYS:HG2	1.47	0.79
1:E:270:HIS:HD2	1:E:283:LYS:HG2	1.47	0.78
1:M:12:VAL:HG12	1:M:12:VAL:O	1.83	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:M:232:PHE:HE1	1:M:354:MET:HE2	1.49	0.78
1:E:212:CYS:HA	8:E:6014:IPA:H33	1.66	0.77
1:I:377:PHE:HB2	1:I:391:VAL:HG22	1.65	0.77
1:E:213:ASP:H	8:E:6014:IPA:H33	1.49	0.77
1:M:291:SER:HB2	10:N:299:HOH:O	1.86	0.76
1:A:27:PRO:HB3	1:A:31:HIS:ND1	2.00	0.76
1:I:253:LEU:CD1	1:I:265:ILE:HD11	2.15	0.76
1:I:12:VAL:HG12	1:I:12:VAL:O	1.86	0.76
1:M:49:ARG:HE	1:M:168:VAL:HG11	1.50	0.76
1:M:253:LEU:CD1	1:M:265:ILE:HD11	2.16	0.76
1:I:4:GLN:HG3	1:I:283:LYS:HE3	1.68	0.75
1:E:9:SER:HB2	1:E:279:THR:OG1	1.87	0.75
1:E:70:VAL:HG21	1:E:251:MET:CE	2.18	0.74
2:N:600:G:O2'	2:N:601:U:H5'	1.87	0.73
1:I:232:PHE:HE1	1:I:354:MET:HE2	1.51	0.72
1:M:225:MET:HE2	1:M:225:MET:HA	1.71	0.72
1:I:49:ARG:HE	1:I:168:VAL:HG11	1.51	0.72
2:J:600:G:O2'	2:J:601:U:H5'	1.89	0.72
1:A:70:VAL:HG21	1:A:251:MET:CE	2.20	0.72
1:M:1:GLY:HA2	1:M:65:ASN:OD1	1.88	0.72
1:A:115:SER:HB2	10:A:545:HOH:O	1.89	0.71
1:M:336:HIS:HB2	8:M:6019:IPA:H31	1.72	0.71
2:F:600:G:O2'	2:F:601:U:H5'	1.90	0.71
1:I:80:HIS:CE1	1:I:318:LEU:HB3	2.25	0.71
1:M:18:ASN:OD1	1:M:276:LYS:HG3	1.91	0.71
1:E:27:PRO:HB3	1:E:31:HIS:ND1	2.04	0.71
1:I:18:ASN:OD1	1:I:276:LYS:HG3	1.90	0.71
1:M:80:HIS:CE1	1:M:318:LEU:HB3	2.24	0.71
1:M:4:GLN:HG3	1:M:283:LYS:HE3	1.71	0.71
1:A:41:ALA:HB2	1:A:163:ARG:HD3	1.73	0.71
1:I:7:ARG:HD2	1:I:12:VAL:CG2	2.21	0.70
1:I:1:GLY:HA2	1:I:65:ASN:OD1	1.91	0.70
1:A:270:HIS:CD2	1:A:283:LYS:HG2	2.27	0.69
1:I:225:MET:HE2	1:I:225:MET:HA	1.73	0.69
1:I:20:PRO:HG3	2:J:598:A:C4	2.27	0.69
1:M:313:TYR:HD2	8:M:6019:IPA:H11	1.57	0.69
1:M:7:ARG:HD2	1:M:12:VAL:CG2	2.22	0.69
1:A:37:VAL:HG13	1:A:165:LYS:HD2	1.75	0.69
1:E:41:ALA:HB2	1:E:163:ARG:HD3	1.75	0.69
1:M:217:PHE:HD1	8:M:6018:IPA:H32	1.56	0.69
2:F:598:A:H3'	2:F:599:G:H5'	1.74	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:381:ASP:HB3	1:E:384:TYR:O	1.92	0.69
2:B:602:C:H2'	2:B:603:U:C6	2.28	0.68
2:B:600:G:O2'	2:B:601:U:H5'	1.92	0.68
1:A:337:GLU:HB2	10:A:571:HOH:O	1.93	0.68
1:E:37:VAL:HG13	1:E:165:LYS:HD2	1.76	0.68
1:I:415:ARG:HH11	1:I:415:ARG:HG2	1.58	0.68
3:C:694:A:N7	10:C:243:HOH:O	2.26	0.68
1:I:253:LEU:HD12	1:I:265:ILE:CD1	2.24	0.68
1:M:311:LYS:O	1:M:311:LYS:HG2	1.93	0.68
2:N:609:C:C4	2:N:610:C:C5	2.82	0.68
1:A:381:ASP:HB3	1:A:384:TYR:O	1.94	0.68
1:M:7:ARG:HH11	1:M:11:GLU:HG2	1.58	0.67
2:B:598:A:H3'	2:B:599:G:H5'	1.75	0.67
1:E:374:LEU:O	1:E:375:LYS:HB2	1.94	0.67
1:E:380:ALA:HA	1:E:388:ILE:HD13	1.78	0.66
2:J:609:C:C4	2:J:610:C:C5	2.84	0.66
1:E:270:HIS:CD2	1:E:283:LYS:HG2	2.27	0.66
1:I:311:LYS:HG2	1:I:311:LYS:O	1.95	0.66
1:E:213:ASP:H	8:E:6014:IPA:C3	2.08	0.66
2:F:602:C:H2'	2:F:603:U:C6	2.30	0.66
1:M:235:THR:HG22	1:M:353:THR:HB	1.78	0.66
1:E:57:ALA:O	1:E:60:SER:HB3	1.96	0.66
1:M:96:CYS:HB3	10:M:536:HOH:O	1.94	0.66
1:A:74:MET:HE1	1:A:244:ALA:HB1	1.78	0.65
1:A:374:LEU:O	1:A:375:LYS:HB2	1.94	0.65
1:E:199:HIS:NE2	10:E:497:HOH:O	2.28	0.65
1:M:20:PRO:HG3	2:N:598:A:C4	2.30	0.65
1:A:380:ALA:HA	1:A:388:ILE:HD13	1.78	0.65
1:A:336:HIS:HB2	8:A:6011:IPA:H11	1.78	0.65
1:A:388:ILE:HD12	8:A:6030:IPA:H32	1.78	0.65
1:M:253:LEU:HD12	1:M:265:ILE:CD1	2.26	0.65
1:M:419:LEU:HD11	2:N:606:C:H4'	1.77	0.64
10:E:560:HOH:O	3:G:697:A:H2	1.80	0.64
1:I:140:GLU:O	1:I:144:LEU:HG	1.97	0.64
1:E:212:CYS:CA	8:E:6014:IPA:H33	2.28	0.64
1:M:446:ASP:C	1:M:447:LEU:HD23	2.23	0.64
1:I:7:ARG:HH11	1:I:11:GLU:HG2	1.60	0.64
3:O:688:G:H8	3:O:688:G:HO5'	1.47	0.63
1:E:213:ASP:N	8:E:6014:IPA:H33	2.13	0.63
1:A:20:PRO:HG3	2:B:598:A:C4	2.34	0.63
1:A:57:ALA:O	1:A:60:SER:HB3	1.98	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:419:LEU:HD11	2:J:606:C:H4'	1.79	0.63
1:I:446:ASP:C	1:I:447:LEU:HD23	2.24	0.63
1:M:49:ARG:HH21	1:M:168:VAL:HG13	1.62	0.63
3:C:688:G:H8	3:C:688:G:HO5'	1.45	0.63
1:E:306:ARG:HG2	1:E:318:LEU:HD13	1.79	0.63
1:I:235:THR:HG22	1:I:353:THR:HB	1.80	0.63
1:A:259:GLY:O	1:A:262:VAL:HG22	1.98	0.63
1:E:21:SER:HB2	1:E:44:THR:HG21	1.81	0.63
1:I:309:LEU:HD12	1:I:321:LEU:HD22	1.81	0.63
1:I:397:ILE:HD13	1:I:421:ALA:HB2	1.80	0.63
1:M:308:LEU:HD12	1:M:354:MET:HE1	1.81	0.62
1:E:74:MET:HE1	1:E:244:ALA:HB1	1.81	0.62
1:E:20:PRO:HG3	2:F:598:A:C4	2.33	0.62
1:I:308:LEU:HD12	1:I:354:MET:HE1	1.80	0.62
1:M:238:ASP:OD1	1:M:239:ALA:N	2.33	0.62
1:M:270:HIS:NE2	1:M:281:CYS:SG	2.73	0.62
1:E:71:ASP:HB3	10:E:482:HOH:O	1.99	0.62
1:M:397:ILE:HD13	1:M:421:ALA:HB2	1.82	0.62
1:A:21:SER:HB2	1:A:44:THR:HG21	1.80	0.61
1:E:47:ASP:OD1	1:E:49:ARG:HD3	1.99	0.61
1:I:49:ARG:HH21	1:I:168:VAL:HG13	1.64	0.61
1:I:356:PRO:HG2	1:I:361:ALA:H	1.66	0.61
1:E:289:GLY:HA2	2:F:600:G:N3	2.16	0.61
1:I:118:TYR:CD1	1:I:119:PRO:HA	2.36	0.61
3:K:688:G:HO5'	3:K:688:G:H8	1.49	0.61
1:E:259:GLY:O	1:E:262:VAL:HG22	2.00	0.61
1:I:199:HIS:NE2	10:I:493:HOH:O	2.31	0.61
1:I:367:THR:HB	10:I:515:HOH:O	2.00	0.61
2:B:598:A:H1'	4:D:805:G:C2	2.36	0.61
1:M:277:ASN:ND2	1:M:278:LYS:HG2	2.15	0.61
1:I:334:TYR:CG	1:I:335:PRO:HD2	2.35	0.60
1:M:356:PRO:HG2	1:M:361:ALA:H	1.66	0.60
1:E:234:TYR:CD1	1:E:328:ASP:HB2	2.36	0.60
1:A:415:ARG:HD3	2:B:606:C:O2'	2.01	0.60
1:M:56:GLU:HG2	10:M:522:HOH:O	2.00	0.60
1:M:140:GLU:O	1:M:144:LEU:HG	2.00	0.60
1:M:309:LEU:HD12	1:M:321:LEU:HD22	1.83	0.60
1:M:234:TYR:CD1	1:M:328:ASP:HB2	2.36	0.60
1:I:270:HIS:CD2	1:I:283:LYS:HE2	2.37	0.60
1:M:270:HIS:CD2	1:M:283:LYS:HE2	2.36	0.60
1:A:58:ILE:HD12	1:A:175:LEU:HD21	1.83	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:37:VAL:O	1:E:37:VAL:HG12	2.02	0.60
1:M:118:TYR:CD1	1:M:119:PRO:HA	2.36	0.60
1:I:238:ASP:OD1	1:I:239:ALA:N	2.35	0.59
1:E:115:SER:HB2	10:E:567:HOH:O	2.02	0.59
3:G:688:G:HO5'	3:G:688:G:H8	1.48	0.59
1:M:415:ARG:HG2	1:M:415:ARG:HH11	1.65	0.59
1:I:277:ASN:ND2	1:I:278:LYS:HG2	2.18	0.59
1:M:334:TYR:CG	1:M:335:PRO:HD2	2.37	0.59
1:A:375:LYS:HD2	3:C:700:G:H5''	1.84	0.59
1:M:6:MET:HG2	1:M:280:TYR:HB3	1.84	0.59
1:M:166:THR:HG22	1:M:170:GLN:NE2	2.18	0.59
1:A:368:TRP:HB3	8:A:6030:IPA:H2	1.84	0.59
1:I:65:ASN:O	1:I:242:SER:HB3	2.02	0.59
1:M:341:SER:HA	1:M:363:PHE:CD2	2.38	0.59
1:I:166:THR:HG22	1:I:170:GLN:NE2	2.17	0.59
1:I:355:THR:HB	1:I:356:PRO:CD	2.32	0.59
1:M:455:ARG:HD3	1:M:455:ARG:O	2.03	0.59
1:E:177:GLU:O	1:E:289:GLY:HA3	2.03	0.59
1:I:341:SER:HA	1:I:363:PHE:CD2	2.38	0.59
1:A:95:MET:O	1:A:189:MET:HG2	2.03	0.58
1:M:235:THR:CG2	1:M:353:THR:HB	2.33	0.58
1:I:426:GLU:N	1:I:450:TYR:CE1	2.71	0.58
1:M:65:ASN:O	1:M:66:LYS:HB2	2.02	0.58
1:M:269:ASN:O	1:M:283:LYS:HA	2.04	0.58
1:M:398:HIS:O	1:M:402:ARG:HG3	2.03	0.58
1:I:6:MET:HG2	1:I:280:TYR:HB3	1.84	0.58
1:A:37:VAL:HG12	1:A:37:VAL:O	2.03	0.58
1:I:235:THR:CG2	1:I:353:THR:HB	2.34	0.58
1:M:355:THR:HB	1:M:356:PRO:CD	2.34	0.58
1:A:306:ARG:HG2	1:A:318:LEU:HD13	1.85	0.58
1:M:313:TYR:CD2	8:M:6019:IPA:H11	2.38	0.58
1:A:234:TYR:CD1	1:A:328:ASP:HB2	2.38	0.58
1:I:8:PRO:HB2	1:I:11:GLU:HB2	1.85	0.58
1:I:234:TYR:CD1	1:I:328:ASP:HB2	2.39	0.58
1:A:397:ILE:HD13	1:A:421:ALA:HB2	1.85	0.57
1:I:232:PHE:HA	1:I:357:ALA:HB2	1.86	0.57
1:I:398:HIS:O	1:I:402:ARG:HG3	2.04	0.57
2:J:605:U:H2'	2:J:606:C:C6	2.38	0.57
1:A:289:GLY:HA2	2:B:600:G:N3	2.19	0.57
1:E:213:ASP:HB3	8:E:6014:IPA:H12	1.86	0.57
1:I:269:ASN:O	1:I:283:LYS:HA	2.04	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:177:GLU:O	1:A:289:GLY:HA3	2.04	0.57
1:I:342:LEU:O	1:I:345:GLN:HB3	2.05	0.57
1:E:425:GLY:HA2	10:E:505:HOH:O	2.03	0.57
2:N:605:U:H2'	2:N:606:C:C6	2.39	0.57
1:E:58:ILE:HD12	1:E:175:LEU:HD21	1.86	0.57
1:E:82:ALA:O	1:E:86:MET:HG2	2.04	0.57
1:I:439:VAL:HG23	1:I:442:GLY:H	1.70	0.57
1:M:11:GLU:O	1:M:12:VAL:HG23	2.05	0.57
1:M:225:MET:HE2	1:M:225:MET:CA	2.34	0.57
1:I:11:GLU:O	1:I:12:VAL:HG23	2.04	0.57
1:M:439:VAL:HG23	1:M:442:GLY:H	1.68	0.57
1:A:82:ALA:O	1:A:86:MET:HG2	2.05	0.57
1:I:133:LYS:HD3	10:I:512:HOH:O	2.04	0.56
1:E:2:GLU:HG3	1:E:64:GLY:HA2	1.87	0.56
1:M:426:GLU:N	1:M:450:TYR:CE1	2.74	0.56
1:A:2:GLU:HG3	1:A:64:GLY:HA2	1.87	0.56
1:A:270:HIS:CD2	1:A:283:LYS:HE2	2.40	0.56
1:E:375:LYS:HD2	3:G:700:G:H5''	1.87	0.56
1:M:49:ARG:HH21	1:M:168:VAL:CG1	2.16	0.56
1:A:419:LEU:HD11	2:B:606:C:H4'	1.87	0.56
1:M:8:PRO:HB2	1:M:11:GLU:HB2	1.86	0.56
1:M:289:GLY:HA2	2:N:600:G:N3	2.20	0.56
1:M:65:ASN:O	1:M:242:SER:HB3	2.05	0.56
1:M:342:LEU:O	1:M:345:GLN:HB3	2.06	0.56
1:E:419:LEU:HD11	2:F:606:C:H4'	1.87	0.56
1:I:270:HIS:CE1	1:I:272:HIS:HE1	2.24	0.56
1:M:356:PRO:HG2	1:M:361:ALA:N	2.21	0.56
1:A:260:ASP:OD1	1:A:260:ASP:N	2.34	0.56
1:E:260:ASP:N	1:E:260:ASP:OD1	2.36	0.56
1:A:52:THR:CG2	1:A:53:ASP:N	2.69	0.55
1:E:397:ILE:HD13	1:E:421:ALA:HB2	1.88	0.55
1:I:426:GLU:HA	1:I:450:TYR:CD1	2.42	0.55
1:A:7:ARG:HD3	1:A:12:VAL:CG2	2.37	0.55
1:E:270:HIS:CD2	1:E:283:LYS:HE2	2.41	0.55
1:I:356:PRO:HG2	1:I:361:ALA:N	2.21	0.55
1:A:112:LEU:HD21	1:A:131:LEU:HB2	1.88	0.55
1:I:455:ARG:HD2	1:I:455:ARG:O	2.05	0.55
1:M:232:PHE:HB3	1:M:355:THR:O	2.07	0.55
1:A:451:SER:O	1:A:455:ARG:HD2	2.07	0.55
1:A:97:LEU:HD13	1:A:189:MET:HE2	1.88	0.55
1:M:49:ARG:HE	1:M:168:VAL:CG1	2.19	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:397:ILE:HD11	1:I:420:LEU:HB3	1.88	0.54
1:A:120:TYR:CE1	1:A:144:LEU:HD13	2.42	0.54
1:I:49:ARG:HH21	1:I:168:VAL:CG1	2.19	0.54
1:I:120:TYR:HB3	1:I:125:LYS:HB3	1.89	0.54
1:E:218:TRP:CD1	1:E:390:PRO:HA	2.43	0.54
1:E:428:GLU:HB2	10:E:579:HOH:O	2.08	0.54
1:I:232:PHE:HB3	1:I:355:THR:O	2.08	0.54
1:M:426:GLU:HA	1:M:450:TYR:CD1	2.43	0.54
1:M:12:VAL:O	1:M:12:VAL:CG1	2.55	0.54
1:E:23:THR:HG22	10:E:547:HOH:O	2.08	0.54
2:F:605:U:H2'	2:F:606:C:C6	2.43	0.54
1:E:112:LEU:HD21	1:E:131:LEU:HB2	1.89	0.53
1:M:120:TYR:HB3	1:M:125:LYS:HB3	1.89	0.53
1:M:133:LYS:HD3	10:M:511:HOH:O	2.08	0.53
1:A:66:LYS:HG2	1:A:349:ASP:O	2.08	0.53
1:I:118:TYR:O	1:I:180:SER:HB2	2.08	0.53
1:E:451:SER:O	1:E:455:ARG:HD2	2.09	0.53
2:J:598:A:H1'	4:L:805:G:N2	2.24	0.53
2:B:602:C:H2'	2:B:603:U:H6	1.73	0.53
2:B:605:U:H2'	2:B:606:C:C6	2.43	0.53
1:I:289:GLY:HA2	2:J:600:G:N3	2.23	0.53
1:M:232:PHE:HA	1:M:357:ALA:HB2	1.89	0.53
1:E:95:MET:O	1:E:189:MET:HG2	2.08	0.53
1:E:97:LEU:HD13	1:E:189:MET:HE2	1.90	0.53
1:I:65:ASN:O	1:I:66:LYS:HB2	2.07	0.53
1:I:374:LEU:O	1:I:375:LYS:HB2	2.08	0.53
1:M:84:GLN:OE1	1:M:306:ARG:NH2	2.41	0.53
1:E:52:THR:CG2	1:E:53:ASP:N	2.72	0.53
1:M:270:HIS:CE1	1:M:272:HIS:HE1	2.27	0.53
1:E:28:SER:HB2	1:E:402:ARG:C	2.33	0.53
1:I:230:PHE:O	1:I:231:ALA:HB2	2.08	0.53
1:I:270:HIS:CE1	1:I:272:HIS:CE1	2.96	0.53
1:M:381:ASP:HB3	1:M:384:TYR:O	2.08	0.53
1:I:84:GLN:OE1	1:I:306:ARG:NH2	2.42	0.53
1:I:225:MET:HE2	1:I:225:MET:CA	2.37	0.53
1:E:70:VAL:CG1	1:E:251:MET:HE1	2.38	0.53
7:I:5002:POP:O4	7:I:5002:POP:O1	2.25	0.53
1:I:74:MET:HE3	1:I:245:TRP:CE3	2.44	0.52
1:E:238:ASP:OD1	1:E:239:ALA:N	2.42	0.52
1:M:277:ASN:HD22	1:M:278:LYS:HG2	1.74	0.52
1:E:70:VAL:HG21	1:E:251:MET:HE3	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:M:234:TYR:HD1	1:M:328:ASP:HB2	1.74	0.52
1:E:225:MET:HA	1:E:225:MET:HE2	1.91	0.52
1:I:277:ASN:HD22	1:I:278:LYS:N	2.08	0.52
1:I:436:ILE:O	1:I:442:GLY:HA3	2.10	0.52
1:E:37:VAL:O	1:E:37:VAL:CG1	2.57	0.52
1:M:74:MET:HE3	1:M:245:TRP:CE3	2.45	0.52
1:I:381:ASP:HB3	1:I:384:TYR:O	2.11	0.51
1:I:320:HIS:HB2	1:I:334:TYR:CE1	2.46	0.51
1:M:277:ASN:HD22	1:M:278:LYS:N	2.08	0.51
2:B:597:C:O2	2:B:597:C:H2'	2.09	0.51
1:E:176:ILE:HG13	2:F:600:G:C8	2.45	0.51
1:E:414:VAL:CG1	1:E:436:ILE:HD13	2.40	0.51
1:I:449:GLU:HB2	10:I:513:HOH:O	2.11	0.51
1:A:238:ASP:OD1	1:A:239:ALA:N	2.43	0.51
2:F:597:C:H2'	2:F:597:C:O2	2.09	0.51
1:I:177:GLU:O	1:I:289:GLY:HA3	2.10	0.51
1:M:177:GLU:O	1:M:289:GLY:HA3	2.11	0.51
1:E:66:LYS:HG2	1:E:349:ASP:O	2.11	0.51
1:M:320:HIS:HB2	1:M:334:TYR:CE1	2.46	0.51
1:I:334:TYR:CD1	1:I:335:PRO:HD2	2.46	0.51
1:M:85:LEU:HA	1:M:88:LEU:HD12	1.93	0.51
1:A:374:LEU:N	10:A:540:HOH:O	2.44	0.50
1:A:37:VAL:O	1:A:37:VAL:CG1	2.60	0.50
1:E:24:LYS:HE3	10:E:524:HOH:O	2.12	0.50
1:I:251:MET:O	1:I:255:LYS:HG3	2.10	0.50
1:M:334:TYR:CD2	1:M:335:PRO:HD2	2.46	0.50
1:I:231:ALA:C	1:I:232:PHE:CD2	2.90	0.50
1:I:232:PHE:CZ	1:I:354:MET:HE2	2.45	0.50
1:M:230:PHE:O	1:M:231:ALA:HB2	2.10	0.50
1:M:270:HIS:CE1	1:M:272:HIS:CE1	2.99	0.50
1:M:358:ASP:O	1:M:359:LYS:HB2	2.11	0.50
1:E:120:TYR:CE1	1:E:144:LEU:HD13	2.46	0.50
1:A:28:SER:HB2	1:A:402:ARG:C	2.37	0.50
1:I:234:TYR:HD1	1:I:328:ASP:HB2	1.77	0.50
1:I:334:TYR:CD2	1:I:335:PRO:HD2	2.47	0.50
1:M:336:HIS:CB	8:M:6019:IPA:H31	2.40	0.50
1:A:27:PRO:HB3	1:A:31:HIS:HD1	1.72	0.50
1:E:39:GLU:OE2	1:E:165:LYS:HG3	2.11	0.50
1:I:334:TYR:CG	1:I:335:PRO:CD	2.94	0.50
1:M:238:ASP:HB2	1:M:286:MET:HB3	1.94	0.50
1:E:236:GLY:HA2	7:G:5003:POP:O6	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:M:231:ALA:C	1:M:232:PHE:CD2	2.89	0.50
2:N:612:G:H5''	2:N:613:A:OP1	2.12	0.50
1:A:313:TYR:CD1	8:A:6011:IPA:H32	2.47	0.50
1:A:316:ILE:HD12	8:A:6011:IPA:H33	1.93	0.50
1:E:414:VAL:HG11	1:E:436:ILE:HD13	1.93	0.50
1:E:446:ASP:C	1:E:447:LEU:HD23	2.37	0.50
1:M:374:LEU:O	1:M:375:LYS:HB2	2.11	0.50
1:I:421:ALA:O	1:I:424:ASN:HB2	2.12	0.49
2:N:609:C:C4	2:N:610:C:H5	2.30	0.49
1:A:40:PRO:HD3	1:A:403:TRP:CH2	2.47	0.49
1:I:79:ASP:OD1	1:I:255:LYS:HE2	2.12	0.49
1:M:218:TRP:CD1	1:M:390:PRO:HA	2.47	0.49
1:E:217:PHE:CD1	8:E:6014:IPA:H31	2.47	0.49
1:M:54:PHE:CE1	1:M:58:ILE:CG2	2.95	0.49
1:M:118:TYR:O	1:M:180:SER:HB2	2.12	0.49
1:M:312:THR:HG22	1:M:313:TYR:CD1	2.48	0.49
1:M:334:TYR:CG	1:M:335:PRO:CD	2.95	0.49
1:E:154:LEU:HD22	1:E:177:GLU:HB3	1.95	0.49
2:J:612:G:H5''	2:J:613:A:OP1	2.13	0.49
1:M:251:MET:O	1:M:255:LYS:HG3	2.11	0.49
1:M:336:HIS:CD2	8:M:6019:IPA:H31	2.48	0.49
1:I:277:ASN:HD22	1:I:277:ASN:C	2.21	0.49
1:M:397:ILE:HD11	1:M:420:LEU:HB3	1.94	0.49
1:A:176:ILE:HG13	2:B:600:G:C8	2.47	0.49
1:A:336:HIS:HB2	8:A:6011:IPA:C1	2.42	0.49
2:B:598:A:C1'	4:D:805:G:N2	2.73	0.49
1:I:277:ASN:HD22	1:I:278:LYS:HG2	1.77	0.49
1:A:70:VAL:CG1	1:A:251:MET:HE1	2.35	0.49
2:F:602:C:H2'	2:F:603:U:H6	1.73	0.49
1:M:277:ASN:HD22	1:M:277:ASN:C	2.21	0.49
1:M:334:TYR:CD1	1:M:335:PRO:HD2	2.48	0.49
1:A:446:ASP:C	1:A:447:LEU:HD23	2.38	0.49
1:E:7:ARG:HD3	1:E:12:VAL:CG2	2.43	0.49
1:I:238:ASP:HB2	1:I:286:MET:HB3	1.95	0.49
2:B:604:C:H2'	2:B:605:U:C6	2.48	0.48
3:K:690:C:H2'	3:K:691:C:O4'	2.13	0.48
2:N:608:U:H2'	2:N:609:C:O4'	2.12	0.48
1:A:81:TYR:O	1:A:84:GLN:HB2	2.13	0.48
1:E:339:ASP:HB3	1:E:342:LEU:HD12	1.95	0.48
2:B:613:A:O5'	3:G:688:G:C5'	2.53	0.48
1:I:358:ASP:O	1:I:359:LYS:HB2	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:154:LEU:HD22	1:A:177:GLU:HB3	1.94	0.48
1:A:384:TYR:CE2	8:A:6028:IPA:H2	2.48	0.48
1:I:53:ASP:HB3	1:I:56:GLU:HB3	1.95	0.48
2:J:609:C:C4	2:J:610:C:H5	2.31	0.48
1:M:97:LEU:O	1:M:101:MET:HG3	2.13	0.48
1:E:216:LEU:HB2	8:E:6014:IPA:H13	1.94	0.48
1:I:9:SER:HA	1:I:279:THR:OG1	2.13	0.48
1:I:54:PHE:CE1	1:I:58:ILE:CG2	2.97	0.48
1:I:128:ARG:HG3	10:I:527:HOH:O	2.14	0.48
1:M:237:TYR:CG	1:M:328:ASP:HB3	2.48	0.48
1:A:146:ASP:HA	10:A:534:HOH:O	2.12	0.48
1:A:358:ASP:OD1	1:A:360:SER:HB3	2.13	0.48
1:I:49:ARG:HE	1:I:168:VAL:CG1	2.21	0.48
1:M:218:TRP:O	1:M:222:PRO:HD3	2.14	0.48
1:I:12:VAL:O	1:I:12:VAL:CG1	2.58	0.48
1:I:97:LEU:O	1:I:101:MET:HG3	2.12	0.48
1:A:232:PHE:HA	1:A:355:THR:O	2.14	0.48
1:I:237:TYR:CG	1:I:328:ASP:HB3	2.49	0.48
1:M:411:GLN:HB3	10:M:496:HOH:O	2.14	0.48
1:I:40:PRO:HD3	1:I:403:TRP:CH2	2.49	0.47
1:I:312:THR:HG22	1:I:313:TYR:CD1	2.49	0.47
2:J:608:U:H2'	2:J:609:C:O4'	2.14	0.47
1:A:45:LYS:H	1:A:45:LYS:HG3	1.39	0.47
1:A:154:LEU:O	1:A:273:HIS:HA	2.14	0.47
1:A:166:THR:O	1:A:170:GLN:HB2	2.13	0.47
1:A:225:MET:HE2	1:A:225:MET:HA	1.96	0.47
1:E:166:THR:O	1:E:170:GLN:HB2	2.15	0.47
2:F:604:C:H2'	2:F:605:U:C6	2.49	0.47
1:I:85:LEU:HA	1:I:88:LEU:HD12	1.95	0.47
1:M:11:GLU:O	1:M:12:VAL:CG2	2.62	0.47
3:O:690:C:H2'	3:O:691:C:O4'	2.14	0.47
1:A:5:TRP:O	1:A:280:TYR:HA	2.14	0.47
1:A:339:ASP:HB3	1:A:342:LEU:HD12	1.95	0.47
1:M:70:VAL:HG21	1:M:251:MET:SD	2.53	0.47
1:M:108:GLU:O	1:M:188:ARG:NH2	2.39	0.47
1:A:39:GLU:OE2	1:A:165:LYS:HG3	2.13	0.47
3:C:690:C:H5''	3:C:690:C:H6	1.80	0.47
1:E:45:LYS:H	1:E:45:LYS:HG3	1.39	0.47
1:I:97:LEU:HD23	1:I:138:THR:HB	1.95	0.47
1:A:387:LEU:HD11	8:A:6028:IPA:H33	1.96	0.47
1:A:398:HIS:O	1:A:402:ARG:HG3	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:H:805:G:H2'	4:H:806:G:O5'	2.15	0.47
1:M:436:ILE:O	1:M:442:GLY:HA3	2.14	0.47
1:E:154:LEU:O	1:E:273:HIS:HA	2.13	0.47
1:M:226:GLU:HG2	1:M:322:LYS:HG3	1.95	0.47
3:C:700:G:H3'	10:C:165:HOH:O	2.14	0.47
1:E:422:TRP:CD1	1:E:453:LEU:HD13	2.50	0.47
3:G:690:C:H5''	3:G:690:C:H6	1.79	0.47
1:I:54:PHE:O	1:I:57:ALA:N	2.45	0.47
1:M:232:PHE:CZ	1:M:354:MET:HE2	2.46	0.47
1:A:270:HIS:NE2	1:A:283:LYS:HE2	2.29	0.47
1:A:414:VAL:CG1	1:A:436:ILE:HD13	2.45	0.47
3:G:688:G:H8	3:G:688:G:O5'	1.98	0.47
1:I:11:GLU:O	1:I:12:VAL:CG2	2.62	0.47
1:M:53:ASP:HB3	1:M:56:GLU:HB3	1.97	0.47
1:A:411:GLN:HE22	1:A:446:ASP:N	2.12	0.47
1:E:81:TYR:O	1:E:84:GLN:HB2	2.14	0.47
1:E:277:ASN:ND2	1:E:278:LYS:HG3	2.30	0.47
1:E:5:TRP:O	1:E:280:TYR:HA	2.15	0.46
1:E:270:HIS:NE2	1:E:283:LYS:HE2	2.30	0.46
1:I:142:GLN:OE1	1:I:142:GLN:HA	2.15	0.46
1:M:40:PRO:HD3	1:M:403:TRP:CH2	2.51	0.46
3:G:692:G:O2'	3:G:693:G:H5'	2.16	0.46
4:H:806:G:O2'	4:H:807:A:H5'	2.15	0.46
1:I:65:ASN:HB3	1:I:243:PRO:HD2	1.97	0.46
1:I:218:TRP:O	1:I:222:PRO:HD3	2.15	0.46
1:A:54:PHE:O	1:A:57:ALA:N	2.45	0.46
1:A:449:GLU:O	1:A:450:TYR:C	2.57	0.46
1:M:421:ALA:O	1:M:424:ASN:HB2	2.15	0.46
1:M:9:SER:HA	1:M:279:THR:OG1	2.15	0.46
1:M:237:TYR:O	1:M:238:ASP:C	2.59	0.46
1:M:253:LEU:CD1	1:M:265:ILE:CD1	2.85	0.46
1:M:339:ASP:OD2	1:M:342:LEU:HG	2.15	0.46
4:D:805:G:H2'	4:D:806:G:O5'	2.15	0.46
1:I:226:GLU:HG2	1:I:322:LYS:HG3	1.97	0.46
1:I:432:PHE:CZ	1:I:436:ILE:HD11	2.50	0.46
4:H:806:G:C2'	4:H:807:A:H5'	2.46	0.46
1:I:102:TYR:CE1	1:I:136:ARG:HA	2.51	0.46
1:M:54:PHE:CE1	1:M:58:ILE:HG21	2.51	0.46
1:E:232:PHE:HA	1:E:355:THR:O	2.16	0.46
1:I:11:GLU:C	1:I:12:VAL:HG23	2.41	0.46
1:I:108:GLU:O	1:I:188:ARG:NH2	2.38	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:M:79:ASP:OD1	1:M:255:LYS:HE2	2.16	0.46
1:A:277:ASN:ND2	1:A:278:LYS:HG3	2.30	0.45
1:E:238:ASP:O	1:E:285:GLY:HA2	2.16	0.45
1:A:334:TYR:CG	1:A:335:PRO:HD2	2.51	0.45
1:I:70:VAL:HG21	1:I:251:MET:SD	2.56	0.45
1:M:25:LEU:HB2	1:M:40:PRO:HG3	1.98	0.45
1:A:338:VAL:HG23	1:A:339:ASP:N	2.31	0.45
1:I:27:PRO:HD3	1:I:403:TRP:CZ3	2.50	0.45
1:I:394:MET:SD	1:I:421:ALA:HB1	2.57	0.45
1:I:337:GLU:OE2	1:I:337:GLU:C	2.60	0.45
1:M:11:GLU:C	1:M:12:VAL:HG23	2.42	0.45
1:E:436:ILE:O	1:E:442:GLY:HA3	2.16	0.45
2:N:609:C:H2'	2:N:610:C:H5'	1.98	0.45
1:A:238:ASP:O	1:A:285:GLY:HA2	2.16	0.45
1:A:431:LYS:HB2	1:A:431:LYS:HE2	1.76	0.45
1:M:97:LEU:HD23	1:M:138:THR:HB	1.99	0.45
1:A:218:TRP:CD1	1:A:390:PRO:HA	2.51	0.45
4:D:806:G:C2'	4:D:807:A:H5'	2.47	0.45
1:E:17:ILE:O	1:E:276:LYS:HA	2.16	0.45
1:E:449:GLU:O	1:E:450:TYR:C	2.59	0.45
1:A:392:MET:HE2	1:A:392:MET:HB2	1.88	0.45
1:A:411:GLN:HE22	1:A:446:ASP:H	1.63	0.45
1:A:102:TYR:CE1	1:A:136:ARG:HA	2.51	0.45
1:I:275:TYR:HD2	1:I:275:TYR:O	1.99	0.45
1:M:339:ASP:CG	1:M:341:SER:HG	2.25	0.45
1:E:334:TYR:CG	1:E:335:PRO:HD2	2.52	0.45
1:M:151:ASN:ND2	1:M:267:TYR:CD2	2.79	0.45
1:A:52:THR:HG23	1:A:53:ASP:N	2.32	0.44
1:E:54:PHE:O	1:E:57:ALA:N	2.46	0.44
1:E:70:VAL:HG21	1:E:251:MET:HE1	1.97	0.44
1:E:329:ASP:OD2	3:G:701:A:H5''	2.18	0.44
2:J:602:C:H2'	2:J:603:U:C6	2.53	0.44
1:M:102:TYR:CE1	1:M:136:ARG:HA	2.52	0.44
1:M:339:ASP:OD2	1:M:341:SER:OG	2.35	0.44
1:A:367:THR:C	1:A:369:GLU:H	2.25	0.44
1:E:40:PRO:HD3	1:E:403:TRP:CH2	2.52	0.44
1:I:218:TRP:CD1	1:I:390:PRO:HA	2.52	0.44
1:A:436:ILE:O	1:A:442:GLY:HA3	2.17	0.44
1:E:358:ASP:OD1	1:E:360:SER:HB3	2.18	0.44
1:E:398:HIS:O	1:E:402:ARG:HG3	2.17	0.44
2:F:598:A:C3'	2:F:599:G:H5'	2.44	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:600:G:C2'	2:F:601:U:H5'	2.48	0.44
1:M:375:LYS:HD3	1:M:396:GLU:OE1	2.18	0.44
1:A:181:LEU:O	1:A:185:VAL:HG23	2.18	0.44
1:E:367:THR:C	1:E:369:GLU:H	2.25	0.44
1:A:395:LYS:HB3	1:A:395:LYS:HE2	1.70	0.44
1:E:394:MET:SD	1:E:421:ALA:HB1	2.57	0.44
1:M:275:TYR:HD2	1:M:275:TYR:O	2.00	0.44
1:M:397:ILE:CD1	1:M:421:ALA:HB2	2.46	0.44
1:E:305:ILE:HD11	1:E:330:VAL:HG11	1.99	0.44
1:I:237:TYR:O	1:I:238:ASP:C	2.59	0.44
1:I:238:ASP:OD1	1:I:238:ASP:C	2.61	0.44
1:I:339:ASP:OD2	1:I:342:LEU:HG	2.18	0.44
1:M:8:PRO:HB2	1:M:11:GLU:CB	2.48	0.44
1:A:364:GLU:HG2	10:A:555:HOH:O	2.18	0.44
1:A:422:TRP:CD1	1:A:453:LEU:HD13	2.52	0.44
1:E:102:TYR:CE1	1:E:136:ARG:HA	2.53	0.44
1:I:440:PRO:HA	1:I:443:ARG:NH1	2.33	0.44
1:M:7:ARG:NH1	1:M:11:GLU:HG2	2.30	0.44
1:M:275:TYR:C	1:M:275:TYR:CD2	2.96	0.44
1:M:276:LYS:HB3	1:M:277:ASN:H	1.60	0.44
1:A:17:ILE:O	1:A:276:LYS:HA	2.17	0.44
3:C:692:G:O2'	3:C:693:G:H5'	2.17	0.44
1:I:339:ASP:OD2	1:I:341:SER:OG	2.35	0.44
3:G:688:G:H1'	10:G:281:HOH:O	2.18	0.43
1:I:8:PRO:HB2	1:I:11:GLU:CB	2.47	0.43
1:I:358:ASP:HB2	1:I:360:SER:OG	2.18	0.43
1:A:414:VAL:HG11	1:A:436:ILE:HD13	2.00	0.43
1:E:338:VAL:HG23	1:E:339:ASP:N	2.33	0.43
1:I:17:ILE:O	1:I:276:LYS:HA	2.18	0.43
1:I:355:THR:CB	1:I:356:PRO:CD	2.96	0.43
1:M:238:ASP:OD1	1:M:238:ASP:C	2.61	0.43
1:A:43:LEU:HD12	1:A:43:LEU:HA	1.75	0.43
2:B:600:G:C2'	2:B:601:U:H5'	2.49	0.43
4:D:806:G:O2'	4:D:807:A:H5'	2.18	0.43
1:E:411:GLN:HE22	1:E:446:ASP:N	2.17	0.43
1:I:54:PHE:CE1	1:I:58:ILE:HG21	2.53	0.43
1:I:222:PRO:HA	1:I:368:TRP:CZ2	2.54	0.43
1:M:2:GLU:HB3	1:M:64:GLY:HA2	2.00	0.43
1:M:228:LYS:O	1:M:333:SER:HB2	2.19	0.43
2:F:609:C:C2'	2:F:610:C:H5'	2.49	0.43
1:A:7:ARG:O	1:A:8:PRO:C	2.61	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:275:TYR:C	1:I:275:TYR:CD2	2.97	0.43
1:M:345:GLN:C	1:M:347:GLY:N	2.76	0.43
1:E:392:MET:HE2	1:E:392:MET:HB2	1.90	0.43
1:I:80:HIS:CE1	1:I:318:LEU:CB	3.01	0.43
1:M:456:ARG:O	1:M:457:TRP:C	2.61	0.43
1:A:49:ARG:HE	1:A:49:ARG:HB2	1.63	0.43
1:A:397:ILE:HG23	1:A:417:LEU:HD13	1.99	0.43
1:I:2:GLU:HB3	1:I:64:GLY:HA2	2.00	0.43
2:J:612:G:C5'	2:J:613:A:OP1	2.67	0.43
1:M:320:HIS:HB3	1:M:335:PRO:CG	2.49	0.43
3:C:688:G:H8	3:C:688:G:O5'	1.98	0.43
3:G:690:C:H6	3:G:690:C:C5'	2.31	0.43
2:J:609:C:H2'	2:J:610:C:H5'	2.00	0.43
1:M:54:PHE:O	1:M:57:ALA:N	2.45	0.43
1:M:222:PRO:HA	1:M:368:TRP:CZ2	2.54	0.43
1:M:276:LYS:HE2	1:M:276:LYS:HB2	1.79	0.43
1:A:27:PRO:HB3	1:A:31:HIS:CE1	2.54	0.42
1:E:4:GLN:HG2	1:E:283:LYS:HE3	2.01	0.42
1:M:17:ILE:O	1:M:276:LYS:HA	2.19	0.42
1:E:43:LEU:HA	1:E:43:LEU:HD12	1.79	0.42
1:M:65:ASN:HB3	1:M:243:PRO:HD2	2.00	0.42
1:M:275:TYR:CE2	1:M:276:LYS:HE2	2.54	0.42
1:M:394:MET:SD	1:M:421:ALA:HB1	2.60	0.42
3:O:694:A:H2'	3:O:695:C:C6	2.54	0.42
3:O:694:A:H2'	3:O:695:C:H6	1.85	0.42
1:I:90:ILE:HD13	1:I:194:LEU:HD12	2.00	0.42
1:I:273:HIS:HB2	1:I:280:TYR:CZ	2.54	0.42
1:A:70:VAL:HG21	1:A:251:MET:HE3	2.00	0.42
1:A:394:MET:SD	1:A:421:ALA:HB1	2.59	0.42
1:E:297:ASN:HB3	1:E:327:GLY:HA2	2.00	0.42
1:E:384:TYR:HA	1:E:385:PRO:HD3	1.83	0.42
1:M:7:ARG:O	1:M:7:ARG:HG3	2.19	0.42
1:M:142:GLN:OE1	1:M:142:GLN:HA	2.19	0.42
1:M:162:LEU:HD11	1:M:403:TRP:CD1	2.54	0.42
2:N:612:G:C5'	2:N:613:A:OP1	2.67	0.42
1:A:422:TRP:CZ2	1:A:423:HIS:CE1	3.07	0.42
1:E:179:SER:HB2	1:E:290:CYS:HA	2.00	0.42
1:E:422:TRP:CZ2	1:E:423:HIS:CE1	3.08	0.42
3:G:701:A:H2'	3:G:702:C:O4'	2.18	0.42
1:I:95:MET:HE3	1:I:95:MET:HB2	1.90	0.42
1:I:154:LEU:HD22	1:I:177:GLU:HB3	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:M:81:TYR:O	1:M:84:GLN:HB2	2.20	0.42
3:C:693:G:H2'	3:C:694:A:H8	1.85	0.42
3:C:701:A:H2'	3:C:702:C:O4'	2.19	0.42
1:E:120:TYR:HB3	1:E:125:LYS:HB3	2.01	0.42
1:E:306:ARG:CG	1:E:318:LEU:HD13	2.46	0.42
1:E:397:ILE:HG23	1:E:417:LEU:HD13	2.02	0.42
1:I:25:LEU:HB2	1:I:40:PRO:HG3	2.01	0.42
1:M:90:ILE:HD13	1:M:194:LEU:HD12	2.01	0.42
1:M:392:MET:HE2	1:M:392:MET:HB2	1.83	0.42
1:A:320:HIS:O	1:A:335:PRO:HD3	2.20	0.42
1:E:95:MET:HE3	1:E:95:MET:HB2	1.86	0.42
1:I:275:TYR:CE2	1:I:276:LYS:HE2	2.53	0.42
1:M:337:GLU:C	1:M:337:GLU:OE2	2.62	0.42
1:A:145:LEU:HA	1:A:145:LEU:HD23	1.87	0.42
1:I:456:ARG:O	1:I:457:TRP:C	2.62	0.42
1:A:269:ASN:ND2	10:A:552:HOH:O	2.49	0.42
1:A:413:HIS:CD2	3:C:698:G:H4'	2.55	0.42
2:J:604:C:H2'	2:J:605:U:O4'	2.19	0.42
1:M:7:ARG:HD3	1:M:11:GLU:HG2	2.02	0.42
2:B:609:C:C2'	2:B:610:C:H5'	2.49	0.41
3:C:690:C:H6	3:C:690:C:C5'	2.32	0.41
1:E:388:ILE:HD12	8:E:6029:IPA:H13	2.03	0.41
1:I:320:HIS:HB3	1:I:335:PRO:CG	2.50	0.41
1:M:118:TYR:CG	1:M:119:PRO:HA	2.55	0.41
1:M:273:HIS:HB2	1:M:280:TYR:CZ	2.55	0.41
1:M:387:LEU:HD22	1:M:461:PHE:CZ	2.55	0.41
1:A:368:TRP:CB	8:A:6030:IPA:H2	2.49	0.41
1:I:397:ILE:CD1	1:I:421:ALA:HB2	2.46	0.41
1:A:35:GLU:O	1:A:402:ARG:HD2	2.20	0.41
2:B:605:U:H5	10:B:492:HOH:O	2.03	0.41
1:I:82:ALA:O	1:I:86:MET:HG2	2.20	0.41
1:M:232:PHE:CB	1:M:355:THR:O	2.69	0.41
1:M:383:LYS:HB3	1:M:384:TYR:CD1	2.56	0.41
1:A:4:GLN:HG2	1:A:283:LYS:HE3	2.02	0.41
1:M:27:PRO:HD3	1:M:403:TRP:CZ3	2.55	0.41
1:M:47:ASP:HA	1:M:48:PRO:HD3	1.89	0.41
1:M:140:GLU:O	1:M:140:GLU:HG3	2.18	0.41
1:M:318:LEU:HA	1:M:318:LEU:HD23	1.74	0.41
1:E:368:TRP:CB	8:E:6029:IPA:H2	2.34	0.41
1:M:355:THR:CB	1:M:356:PRO:CD	2.99	0.41
1:A:97:LEU:O	1:A:101:MET:HG3	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:218:TRP:O	1:A:222:PRO:HD3	2.21	0.41
3:G:693:G:H2'	3:G:694:A:H8	1.85	0.41
1:M:447:LEU:HD23	1:M:447:LEU:N	2.34	0.41
1:A:306:ARG:CG	1:A:318:LEU:HD13	2.51	0.41
1:A:384:TYR:HA	1:A:385:PRO:HD3	1.82	0.41
2:B:598:A:C3'	2:B:599:G:H5'	2.46	0.41
1:E:50:LEU:HD21	1:E:171:GLY:HA3	2.03	0.41
1:I:7:ARG:HD3	1:I:11:GLU:HG2	2.02	0.41
1:I:118:TYR:CG	1:I:119:PRO:HA	2.55	0.41
1:A:55:GLU:O	1:A:59:PHE:HD2	2.03	0.41
1:A:381:ASP:OD1	8:A:6027:IPA:H2	2.21	0.41
1:M:321:LEU:HD12	1:M:322:LYS:N	2.35	0.41
1:A:24:LYS:HB2	1:A:24:LYS:HE3	1.76	0.41
1:A:217:PHE:HD1	8:A:6009:IPA:H12	1.85	0.41
1:A:305:ILE:HD11	1:A:330:VAL:HG11	2.03	0.41
1:E:27:PRO:HB3	1:E:31:HIS:CE1	2.55	0.41
1:E:52:THR:HG23	1:E:53:ASP:N	2.35	0.41
1:E:226:GLU:HG3	1:E:320:HIS:O	2.21	0.41
1:E:295:ILE:HG12	10:E:507:HOH:O	2.21	0.41
1:E:411:GLN:HE22	1:E:446:ASP:H	1.69	0.41
1:I:228:LYS:O	1:I:333:SER:HB2	2.21	0.41
1:I:275:TYR:O	1:I:275:TYR:CD2	2.74	0.41
2:J:604:C:H2'	2:J:605:U:C6	2.56	0.41
3:K:694:A:H2'	3:K:695:C:C6	2.56	0.41
1:M:82:ALA:O	1:M:86:MET:HG2	2.21	0.41
1:M:275:TYR:O	1:M:276:LYS:CB	2.69	0.41
2:N:604:C:H2'	2:N:605:U:C6	2.55	0.41
3:O:692:G:O2'	3:O:693:G:H5'	2.21	0.41
1:A:120:TYR:HB3	1:A:125:LYS:HB3	2.03	0.41
2:B:610:C:C2'	2:B:611:G:O5'	2.69	0.41
1:E:213:ASP:HB3	8:E:6014:IPA:C1	2.50	0.41
1:E:254:GLU:HG2	1:E:259:GLY:HA2	2.01	0.41
1:I:275:TYR:O	1:I:276:LYS:CB	2.69	0.41
1:M:432:PHE:CZ	1:M:436:ILE:HD11	2.55	0.41
2:N:604:C:H2'	2:N:605:U:O4'	2.21	0.41
2:F:598:A:H1'	4:H:805:G:C2	2.57	0.40
1:I:47:ASP:HA	1:I:48:PRO:HD3	1.87	0.40
1:I:270:HIS:NE2	1:I:283:LYS:HE2	2.36	0.40
1:M:226:GLU:HG3	1:M:320:HIS:O	2.20	0.40
1:A:179:SER:HB2	1:A:290:CYS:HA	2.02	0.40
1:A:254:GLU:HG2	1:A:259:GLY:HA2	2.02	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:382:GLU:HB2	8:A:6027:IPA:H33	2.03	0.40
1:E:212:CYS:HB2	8:E:6014:IPA:C3	2.51	0.40
1:E:237:TYR:CG	1:E:328:ASP:HB3	2.56	0.40
2:F:610:C:C2'	2:F:611:G:O5'	2.69	0.40
1:I:162:LEU:HD11	1:I:403:TRP:CD1	2.56	0.40
1:I:318:LEU:HD23	1:I:318:LEU:HA	1.75	0.40
1:M:448:PRO:HG2	1:M:453:LEU:HD21	2.02	0.40
1:A:20:PRO:HA	2:B:598:A:N1	2.36	0.40
1:E:223:VAL:O	1:E:223:VAL:HG23	2.21	0.40
1:E:334:TYR:CD2	1:E:335:PRO:HD2	2.55	0.40
1:I:375:LYS:HD3	1:I:396:GLU:OE1	2.21	0.40
1:I:457:TRP:O	1:I:458:LEU:C	2.64	0.40
1:M:67:ILE:HG23	1:M:244:ALA:HB3	2.04	0.40
1:M:457:TRP:O	1:M:458:LEU:C	2.63	0.40
1:A:58:ILE:HD11	1:A:59:PHE:CE2	2.57	0.40
1:A:214:PRO:O	1:A:390:PRO:HG3	2.22	0.40
1:E:61:LYS:HD2	1:E:239:ALA:CB	2.51	0.40
1:E:121:VAL:CG1	1:E:122:ALA:N	2.85	0.40
3:G:693:G:H2'	3:G:694:A:C8	2.57	0.40
1:M:270:HIS:NE2	1:M:283:LYS:HE2	2.36	0.40
1:M:345:GLN:C	1:M:347:GLY:H	2.27	0.40
1:E:145:LEU:HD23	1:E:145:LEU:HA	1.90	0.40
1:I:275:TYR:O	1:I:276:LYS:HB3	2.22	0.40
1:I:321:LEU:HD12	1:I:322:LYS:N	2.36	0.40
1:I:334:TYR:CD1	1:I:335:PRO:CD	3.05	0.40
1:I:355:THR:HB	1:I:356:PRO:HD3	2.03	0.40
1:I:383:LYS:HB3	1:I:384:TYR:CD1	2.57	0.40
1:I:411:GLN:HB3	10:I:546:HOH:O	2.22	0.40
1:M:334:TYR:CD1	1:M:335:PRO:CD	3.04	0.40
1:M:358:ASP:HB2	1:M:360:SER:OG	2.22	0.40
1:M:440:PRO:HA	1:M:443:ARG:NH1	2.36	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	459/471 (98%)	434 (95%)	23 (5%)	2 (0%)	30	50
1	E	459/471 (98%)	432 (94%)	25 (5%)	2 (0%)	30	50
1	I	459/471 (98%)	406 (88%)	45 (10%)	8 (2%)	7	14
1	M	459/471 (98%)	407 (89%)	44 (10%)	8 (2%)	7	14
All	All	1836/1884 (98%)	1679 (91%)	137 (8%)	20 (1%)	11	22

All (20) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	285	GLY
1	E	285	GLY
1	I	12	VAL
1	I	64	GLY
1	I	70	VAL
1	I	164	SER
1	M	64	GLY
1	M	70	VAL
1	M	164	SER
1	I	66	LYS
1	M	12	VAL
1	M	66	LYS
1	I	276	LYS
1	I	355	THR
1	I	457	TRP
1	M	276	LYS
1	M	355	THR
1	M	457	TRP
1	E	70	VAL
1	A	70	VAL

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	403/412 (98%)	365 (91%)	38 (9%)	8	16
1	E	403/412 (98%)	365 (91%)	38 (9%)	8	16
1	I	403/412 (98%)	364 (90%)	39 (10%)	8	15
1	M	403/412 (98%)	365 (91%)	38 (9%)	8	16
All	All	1612/1648 (98%)	1459 (90%)	153 (10%)	8	16

All (153) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	GLU
1	A	4	GLN
1	A	6	MET
1	A	9	SER
1	A	10	LYS
1	A	17	ILE
1	A	24	LYS
1	A	28	SER
1	A	35	GLU
1	A	43	LEU
1	A	45	LYS
1	A	49	ARG
1	A	52	THR
1	A	94	GLN
1	A	139	LYS
1	A	163	ARG
1	A	166	THR
1	A	170	GLN
1	A	176	ILE
1	A	200	LYS
1	A	220	LYS
1	A	223	VAL
1	A	260	ASP
1	A	261	ARG
1	A	294	SER
1	A	314	LYS
1	A	316	ILE
1	A	329	ASP
1	A	330	VAL
1	A	337	GLU
1	A	338	VAL
1	A	348	LYS
1	A	359	LYS

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Mol	Chain	Res	Type
1	A	392	MET
1	A	415	ARG
1	A	428	GLU
1	A	455	ARG
1	A	456	ARG
1	E	2	GLU
1	E	4	GLN
1	E	6	MET
1	E	9	SER
1	E	10	LYS
1	E	17	ILE
1	E	24	LYS
1	E	28	SER
1	E	35	GLU
1	E	43	LEU
1	E	45	LYS
1	E	49	ARG
1	E	52	THR
1	E	84	GLN
1	E	94	GLN
1	E	139	LYS
1	E	163	ARG
1	E	166	THR
1	E	170	GLN
1	E	176	ILE
1	E	200	LYS
1	E	220	LYS
1	E	223	VAL
1	E	260	ASP
1	E	261	ARG
1	E	294	SER
1	E	314	LYS
1	E	316	ILE
1	E	329	ASP
1	E	330	VAL
1	E	337	GLU
1	E	338	VAL
1	E	348	LYS
1	E	359	LYS
1	E	392	MET
1	E	428	GLU
1	E	455	ARG

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Mol	Chain	Res	Type
1	E	456	ARG
1	I	3	ILE
1	I	5	TRP
1	I	6	MET
1	I	21	SER
1	I	26	GLU
1	I	39	GLU
1	I	56	GLU
1	I	60	SER
1	I	63	VAL
1	I	68	THR
1	I	84	GLN
1	I	121	VAL
1	I	151	ASN
1	I	166	THR
1	I	176	ILE
1	I	179	SER
1	I	180	SER
1	I	261	ARG
1	I	277	ASN
1	I	291	SER
1	I	294	SER
1	I	295	ILE
1	I	316	ILE
1	I	329	ASP
1	I	330	VAL
1	I	331	ILE
1	I	337	GLU
1	I	338	VAL
1	I	346	SER
1	I	348	LYS
1	I	352	LEU
1	I	364	GLU
1	I	367	THR
1	I	391	VAL
1	I	415	ARG
1	I	428	GLU
1	I	431	LYS
1	I	441	ILE
1	I	455	ARG
1	M	3	ILE
1	M	5	TRP

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Mol	Chain	Res	Type
1	M	6	MET
1	M	21	SER
1	M	26	GLU
1	M	39	GLU
1	M	56	GLU
1	M	60	SER
1	M	63	VAL
1	M	68	THR
1	M	84	GLN
1	M	121	VAL
1	M	151	ASN
1	M	166	THR
1	M	176	ILE
1	M	179	SER
1	M	180	SER
1	M	261	ARG
1	M	277	ASN
1	M	291	SER
1	M	294	SER
1	M	295	ILE
1	M	316	ILE
1	M	329	ASP
1	M	330	VAL
1	M	337	GLU
1	M	338	VAL
1	M	346	SER
1	M	348	LYS
1	M	352	LEU
1	M	364	GLU
1	M	367	THR
1	M	391	VAL
1	M	415	ARG
1	M	428	GLU
1	M	431	LYS
1	M	441	ILE
1	M	455	ARG

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

Mol	Chain	Res	Type
1	A	18	ASN
1	A	94	GLN

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Mol	Chain	Res	Type
1	A	170	GLN
1	A	277	ASN
1	A	336	HIS
1	A	411	GLN
1	E	18	ASN
1	E	94	GLN
1	E	170	GLN
1	E	277	ASN
1	I	170	GLN
1	I	269	ASN
1	I	270	HIS
1	I	272	HIS
1	I	277	ASN
1	I	398	HIS
1	M	4	GLN
1	M	170	GLN
1	M	269	ASN
1	M	270	HIS
1	M	272	HIS
1	M	277	ASN
1	M	336	HIS
1	M	398	HIS

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	B	15/26 (57%)	3 (20%)	2 (13%)
2	F	15/26 (57%)	3 (20%)	2 (13%)
2	J	15/26 (57%)	2 (13%)	1 (6%)
2	N	15/26 (57%)	2 (13%)	1 (6%)
3	C	14/15 (93%)	1 (7%)	0
3	G	14/15 (93%)	1 (7%)	0
3	K	14/15 (93%)	0	0
3	O	14/15 (93%)	0	0
4	D	2/9 (22%)	2 (100%)	0
4	H	2/9 (22%)	2 (100%)	0
4	L	3/9 (33%)	1 (33%)	0
4	P	3/9 (33%)	1 (33%)	0
All	All	126/200 (63%)	18 (14%)	6 (4%)

All (18) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
2	B	598	A
2	B	599	G
2	B	600	G
3	C	690	C
4	D	806	G
4	D	807	A
2	F	598	A
2	F	599	G
2	F	600	G
3	G	690	C
4	H	806	G
4	H	807	A
2	J	611	G
2	J	612	G
4	L	807	A
2	N	611	G
2	N	612	G
4	P	807	A

All (6) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
2	B	598	A
2	B	599	G
2	F	598	A
2	F	599	G
2	J	598	A
2	N	598	A

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 29 ligands modelled in this entry, 12 are monoatomic - leaving 17 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	$\# Z > 2$
8	IPA	A	6030	-	3,3,3	0.51	0	3,3,3	0.51	0
8	IPA	M	6018	-	3,3,3	0.63	0	3,3,3	0.25	0
8	IPA	E	6029	-	3,3,3	0.54	0	3,3,3	0.23	0
7	POP	A	5004	6	6,8,8	0.75	0	12,13,13	1.19	1 (8%)
7	POP	G	5003	6	6,8,8	0.69	0	12,13,13	0.97	1 (8%)
7	POP	O	5001	6	6,8,8	0.63	0	12,13,13	1.09	1 (8%)
7	POP	I	5002	6	6,8,8	0.71	0	12,13,13	0.94	1 (8%)
8	IPA	A	6028	-	3,3,3	0.66	0	3,3,3	0.27	0
9	GOL	A	8012	-	5,5,5	0.39	0	5,5,5	0.49	0
9	GOL	J	8005	-	5,5,5	0.43	0	5,5,5	0.94	0
9	GOL	N	8010	-	5,5,5	0.39	0	5,5,5	0.98	0
8	IPA	A	6009	-	3,3,3	0.66	0	3,3,3	0.21	0
9	GOL	E	8011	5	5,5,5	0.33	0	5,5,5	0.45	0
8	IPA	M	6019	-	3,3,3	0.52	0	3,3,3	0.40	0
8	IPA	A	6027	-	3,3,3	0.62	0	3,3,3	0.15	0
8	IPA	A	6011	-	3,3,3	0.51	0	3,3,3	0.38	0
8	IPA	E	6014	-	3,3,3	0.41	0	3,3,3	0.66	0

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
7	POP	A	5004	6	-	3/6/6/6	-
7	POP	G	5003	6	-	1/6/6/6	-
7	POP	O	5001	6	-	0/6/6/6	-
7	POP	I	5002	6	-	0/6/6/6	-
9	GOL	A	8012	-	-	0/4/4/4	-
9	GOL	J	8005	-	-	2/4/4/4	-
9	GOL	N	8010	-	-	4/4/4/4	-
9	GOL	E	8011	5	-	2/4/4/4	-

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
7	O	5001	POP	O2-P1-O	2.73	113.79	104.64
7	G	5003	POP	O6-P2-O	2.72	113.75	104.64
7	A	5004	POP	O5-P2-O	2.45	112.84	104.64
7	I	5002	POP	O3-P1-O	2.11	111.71	104.64

There are no chirality outliers.

All (12) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	A	5004	POP	P2-O-P1-O2
9	N	8010	GOL	O1-C1-C2-C3
9	E	8011	GOL	C1-C2-C3-O3
9	J	8005	GOL	O1-C1-C2-C3
9	N	8010	GOL	C1-C2-C3-O3
9	N	8010	GOL	O1-C1-C2-O2
7	A	5004	POP	P1-O-P2-O4
7	G	5003	POP	P2-O-P1-O2
9	E	8011	GOL	O2-C2-C3-O3
9	J	8005	GOL	O1-C1-C2-O2
7	A	5004	POP	P2-O-P1-O1
9	N	8010	GOL	O2-C2-C3-O3

There are no ring outliers.

11 monomers are involved in 34 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
8	A	6030	IPA	3	0
8	M	6018	IPA	1	0
8	E	6029	IPA	3	0
7	G	5003	POP	1	0
7	I	5002	POP	1	0
8	A	6028	IPA	2	0
8	A	6009	IPA	1	0
8	M	6019	IPA	5	0
8	A	6027	IPA	2	0
8	A	6011	IPA	4	0
8	E	6014	IPA	11	0

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	461/471 (97%)	0.71	38 (8%) 17 18	37, 61, 92, 111	0
1	E	461/471 (97%)	0.62	37 (8%) 18 18	39, 60, 93, 110	0
1	I	461/471 (97%)	0.89	52 (11%) 10 8	40, 68, 115, 143	0
1	M	461/471 (97%)	0.88	62 (13%) 7 5	38, 68, 115, 143	0
2	B	17/26 (65%)	0.22	2 (11%) 9 7	47, 62, 135, 164	0
2	F	17/26 (65%)	0.78	3 (17%) 4 3	47, 62, 134, 164	0
2	J	17/26 (65%)	0.70	3 (17%) 4 3	53, 75, 166, 182	0
2	N	17/26 (65%)	1.37	4 (23%) 2 2	51, 76, 166, 182	0
3	C	15/15 (100%)	0.75	3 (20%) 3 2	49, 65, 124, 141	0
3	G	15/15 (100%)	0.91	2 (13%) 7 6	49, 65, 124, 141	0
3	K	15/15 (100%)	1.12	2 (13%) 7 6	52, 83, 145, 149	0
3	O	15/15 (100%)	1.09	3 (20%) 3 2	52, 84, 145, 149	0
4	D	3/9 (33%)	1.92	1 (33%) 1 0	168, 168, 172, 173	0
4	H	3/9 (33%)	2.85	2 (66%) 0 0	168, 168, 172, 173	0
4	L	4/9 (44%)	1.43	0 100 100	146, 152, 157, 173	0
4	P	4/9 (44%)	1.75	1 (25%) 2 1	146, 152, 158, 173	0
All	All	1986/2084 (95%)	0.79	215 (10%) 11 9	37, 64, 114, 182	0

All (215) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	284	GLY	7.3
1	E	285	GLY	6.4
1	I	357	ALA	6.1
1	A	184	SER	5.4
1	A	64	GLY	5.2

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Mol	Chain	Res	Type	RSRZ
1	M	460	SER	5.0
1	M	357	ALA	4.7
1	M	134	GLN	4.7
1	E	283	LYS	4.6
1	M	287	PRO	4.5
1	M	137	ASP	4.3
1	A	113	SER	4.2
1	I	134	GLN	4.1
1	I	412	ASP	4.1
1	A	357	ALA	4.0
1	A	256	ILE	4.0
1	A	456	ARG	4.0
1	M	71	ASP	4.0
3	C	688	G	3.8
1	A	19	ALA	3.8
1	M	9	SER	3.8
4	H	805	G	3.7
1	E	233	ASP	3.7
1	M	284	GLY	3.7
1	M	290	CYS	3.7
3	K	689	C	3.6
1	M	285	GLY	3.6
2	B	597	C	3.5
1	M	178	ALA	3.5
1	M	8	PRO	3.5
1	M	38	LYS	3.5
1	A	134	GLN	3.5
3	K	688	G	3.5
1	M	67	ILE	3.5
1	I	151	ASN	3.5
1	M	10	LYS	3.5
2	F	597	C	3.4
1	E	137	ASP	3.4
2	N	613	A	3.4
1	E	134	GLN	3.4
1	I	320	HIS	3.4
1	E	456	ARG	3.4
1	E	178	ALA	3.3
1	M	192	GLY	3.3
1	I	46	ASN	3.3
1	I	89	ASP	3.3
1	A	291	SER	3.3

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Mol	Chain	Res	Type	RSRZ
3	O	688	G	3.3
1	E	2	GLU	3.3
1	M	204	VAL	3.3
1	I	313	TYR	3.2
1	I	256	ILE	3.2
1	I	285	GLY	3.2
1	M	164	SER	3.2
1	A	12	VAL	3.2
2	N	597	C	3.1
1	M	277	ASN	3.1
1	M	358	ASP	3.1
3	G	688	G	3.1
1	A	8	PRO	3.1
1	M	70	VAL	3.0
1	M	356	PRO	3.0
1	M	342	LEU	3.0
1	E	138	THR	3.0
1	M	343	LEU	2.9
1	A	359	LYS	2.9
2	N	612	G	2.9
1	A	32	TYR	2.9
4	H	806	G	2.9
1	E	317	ASP	2.9
1	I	1	GLY	2.9
1	M	312	THR	2.9
1	I	310	LEU	2.9
1	M	16	ILE	2.9
1	I	342	LEU	2.9
1	I	67	ILE	2.8
1	M	254	GLU	2.8
1	M	396	GLU	2.8
1	A	414	VAL	2.8
3	G	689	C	2.8
1	A	67	ILE	2.8
1	I	343	LEU	2.8
1	M	339	ASP	2.7
1	M	32	TYR	2.7
2	J	612	G	2.7
1	M	245	TRP	2.7
1	A	100	ALA	2.7
1	M	345	GLN	2.7
1	M	236	GLY	2.7

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Mol	Chain	Res	Type	RSRZ
1	M	5	TRP	2.7
1	E	282	VAL	2.7
1	M	44	THR	2.7
2	B	598	A	2.6
1	I	361	ALA	2.6
1	I	32	TYR	2.6
1	E	290	CYS	2.6
1	E	72	GLU	2.6
1	M	2	GLU	2.6
1	M	138	THR	2.6
1	A	154	LEU	2.6
1	M	415	ARG	2.6
1	M	172	LYS	2.6
3	O	696	G	2.6
1	I	356	PRO	2.6
1	I	358	ASP	2.6
1	M	359	LYS	2.6
1	E	19	ALA	2.6
1	I	47	ASP	2.5
1	I	443	ARG	2.5
1	A	313	TYR	2.5
1	A	33	VAL	2.5
1	M	235	THR	2.5
1	I	327	GLY	2.5
1	M	286	MET	2.5
3	C	689	C	2.5
1	M	65	ASN	2.5
1	M	360	SER	2.5
1	A	455	ARG	2.5
1	E	71	ASP	2.5
1	I	10	LYS	2.5
1	I	449	GLU	2.5
1	A	285	GLY	2.5
1	I	171	GLY	2.5
4	D	805	G	2.5
1	I	9	SER	2.5
1	I	2	GLU	2.5
1	E	287	PRO	2.5
1	E	49	ARG	2.4
1	I	334	TYR	2.4
1	A	276	LYS	2.4
1	E	170	GLN	2.4

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Mol	Chain	Res	Type	RSRZ
1	I	414	VAL	2.4
1	I	461	PHE	2.4
1	M	456	ARG	2.4
1	I	249	LEU	2.4
1	E	62	TYR	2.4
1	A	44	THR	2.4
1	M	355	THR	2.4
1	I	460	SER	2.4
1	E	289	GLY	2.4
1	I	448	PRO	2.4
1	E	4	GLN	2.4
1	E	316	ILE	2.4
1	I	245	TRP	2.4
1	A	1	GLY	2.4
1	A	296	PHE	2.4
1	A	396	GLU	2.3
3	C	690	C	2.3
1	A	345	GLN	2.3
1	I	345	GLN	2.3
1	E	313	TYR	2.3
1	E	179	SER	2.3
1	I	113	SER	2.3
1	M	351	GLY	2.3
2	J	597	C	2.3
1	M	17	ILE	2.3
1	A	18	ASN	2.3
1	A	21	SER	2.3
1	M	315	GLY	2.3
1	E	59	PHE	2.3
1	E	67	ILE	2.3
4	P	805	G	2.3
1	I	65	ASN	2.3
1	A	242	SER	2.3
1	M	63	VAL	2.3
1	I	173	SER	2.2
1	M	446	ASP	2.2
1	A	249	LEU	2.2
1	E	16	ILE	2.2
1	M	363	PHE	2.2
1	I	100	ALA	2.2
2	F	598	A	2.2
1	I	15	PRO	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	70	VAL	2.2
1	M	338	VAL	2.2
1	I	23	THR	2.2
1	I	118	TYR	2.2
1	A	233	ASP	2.2
1	I	438	SER	2.2
1	E	232	PHE	2.2
2	F	603	U	2.2
1	E	189	MET	2.2
3	O	689	C	2.1
1	E	10	LYS	2.1
1	E	275	TYR	2.1
1	M	170	GLN	2.1
1	M	43	LEU	2.1
1	A	42	VAL	2.1
1	E	44	THR	2.1
1	A	175	LEU	2.1
1	I	238	ASP	2.1
1	I	441	ILE	2.1
1	M	215	ASP	2.1
2	J	598	A	2.1
1	I	353	THR	2.1
1	I	359	LYS	2.1
1	A	103	GLY	2.1
1	E	13	GLY	2.1
1	I	103	GLY	2.1
1	A	215	ASP	2.1
1	E	405	LYS	2.1
1	M	362	THR	2.1
1	I	110	LEU	2.1
1	I	61	LYS	2.0
1	M	23	THR	2.0
1	I	286	MET	2.0
2	N	598	A	2.0
1	M	18	ASN	2.0
1	E	17	ILE	2.0
1	A	96	CYS	2.0
1	M	313	TYR	2.0
1	E	286	MET	2.0
1	I	457	TRP	2.0
1	M	352	LEU	2.0
1	M	336	HIS	2.0

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

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

6.3 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
8	IPA	A	6027	4/4	0.52	0.35	84,91,95,98	0
9	GOL	A	8012	6/6	0.58	0.22	71,86,94,98	0
8	IPA	A	6028	4/4	0.61	0.20	60,80,87,112	0
7	POP	I	5002	9/9	0.65	0.22	97,119,129,136	9
7	POP	G	5003	9/9	0.66	0.22	47,60,88,89	9
9	GOL	E	8011	6/6	0.70	0.16	69,82,89,94	0
8	IPA	M	6019	4/4	0.76	0.23	80,96,100,105	0
8	IPA	A	6011	4/4	0.77	0.25	65,88,92,106	0
7	POP	A	5004	9/9	0.78	0.16	45,77,111,115	9
8	IPA	A	6009	4/4	0.80	0.25	64,72,75,78	0
7	POP	O	5001	9/9	0.82	0.13	41,73,93,98	9
9	GOL	J	8005	6/6	0.84	0.15	47,55,64,73	0
8	IPA	M	6018	4/4	0.85	0.14	47,52,58,72	0
8	IPA	A	6030	4/4	0.88	0.19	68,73,76,78	0
9	GOL	N	8010	6/6	0.89	0.24	55,71,72,86	0
5	ZN	I	2002	1/1	0.91	0.08	70,70,70,70	1
8	IPA	E	6029	4/4	0.91	0.15	67,75,80,84	0
5	ZN	M	2001	1/1	0.92	0.10	70,70,70,70	1
8	IPA	E	6014	4/4	0.94	0.16	53,58,75,75	0
6	MN	M	3001	1/1	0.95	0.10	127,127,127,127	0
5	ZN	A	2004	1/1	0.95	0.06	81,81,81,81	1
6	MN	I	3004	1/1	0.95	0.06	90,90,90,90	0
6	MN	M	3002	1/1	0.96	0.06	86,86,86,86	0
6	MN	E	3005	1/1	0.96	0.05	87,87,87,87	0
6	MN	I	3003	1/1	0.97	0.05	120,120,120,120	0
6	MN	A	3007	1/1	0.97	0.04	91,91,91,91	0
5	ZN	E	2003	1/1	0.97	0.09	88,88,88,88	1

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
6	MN	E	3006	1/1	0.98	0.04	74,74,74,74	0
6	MN	A	3008	1/1	0.98	0.04	74,74,74,74	0

6.5 Other polymers [i](#)

There are no such residues in this entry.