



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 12:23 PM UTC

PDB ID : 3OLB / pdb_00003olb
Title : Poliovirus polymerase elongation complex with 2',3'-dideoxy-ctp
Authors : Gong, P.; Peersen, O.B.
Deposited on : 2010-08-25
Resolution : 2.41 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
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

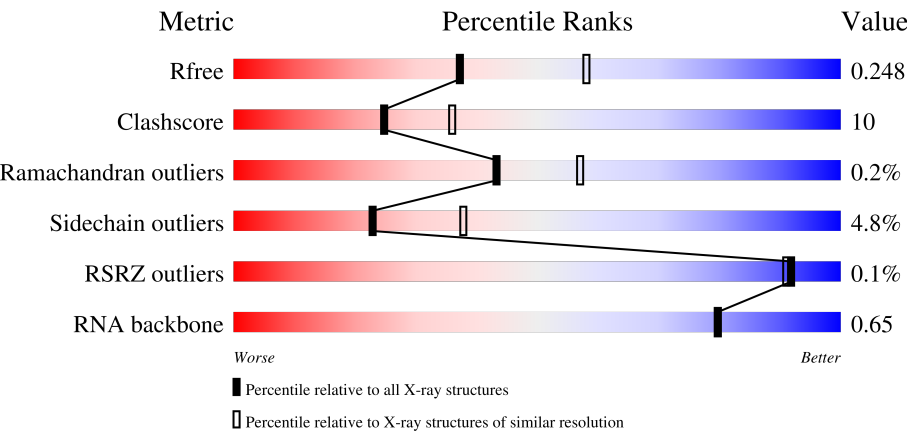
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.41 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.



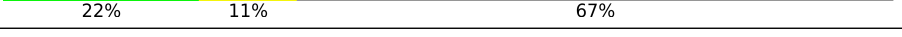
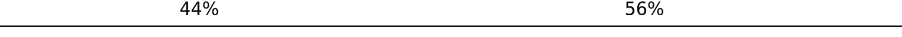
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R _{free}	180053	6062 (2.44-2.40)
Clashscore	190562	6562 (2.44-2.40)
Ramachandran outliers	187476	6481 (2.44-2.40)
Sidechain outliers	187428	6482 (2.44-2.40)
RSRZ outliers	180081	6066 (2.44-2.40)
RNA backbone	3983	1058 (2.70-2.14)

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	78% 18% ..
1	E	471	77% 18% ..
1	I	471	81% 16% ..
1	M	471	81% 15% ..

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Mol	Chain	Length	Quality of chain
2	B	26	
2	F	26	
2	J	26	
2	N	26	
3	C	14	
3	G	14	
3	K	14	
3	O	14	
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
7	IPA	A	6032	-	-	X	-
7	IPA	M	6025	-	-	X	-
7	IPA	M	6031	-	-	X	-

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 18860 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*U*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	18	Total	C	N	O	P	0	0	0
			361	159	59	125	18			
2	F	18	Total	C	N	O	P	0	0	0
			361	159	59	125	18			
2	J	19	Total	C	N	O	P	0	0	0
			381	168	61	133	19			
2	N	19	Total	C	N	O	P	0	0	0
			381	168	61	133	19			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	14	Total 303	C 136	N 62	O 92	P 13	0	0	0
3	G	14	Total 303	C 136	N 62	O 92	P 13	0	0	0
3	K	14	Total 303	C 136	N 62	O 92	P 13	0	0	0
3	O	14	Total 303	C 136	N 62	O 92	P 13	0	0	0

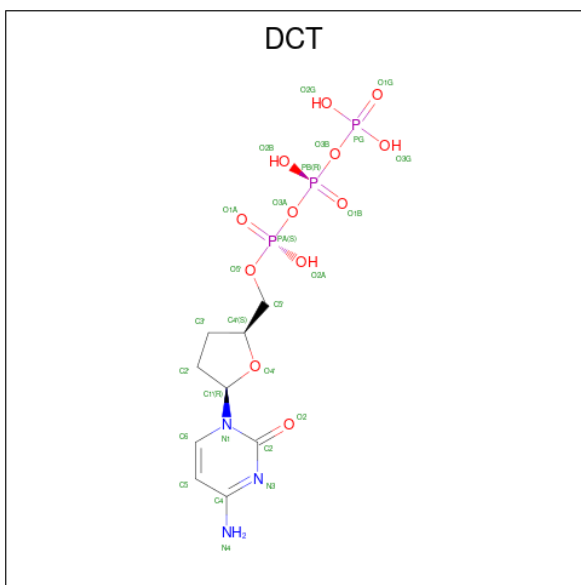
- Molecule 4 is a RNA chain called RNA (5'-R(*GP*GP*GP*AP*GP*AP*UP*GP*A)-3').

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

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

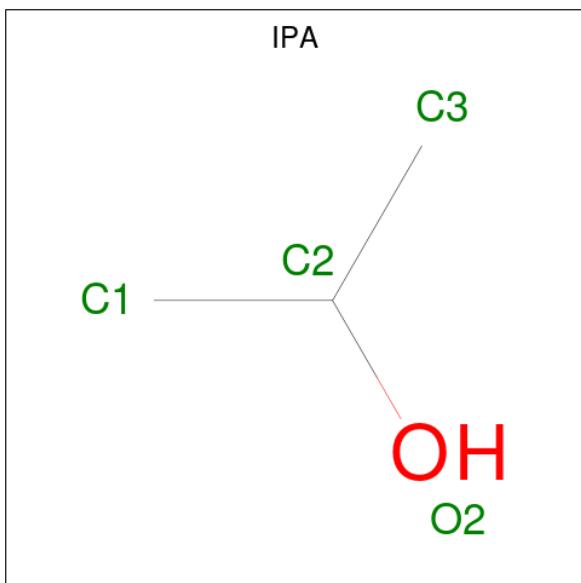
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total 1	Zn 1	0	0
5	E	1	Total 1	Zn 1	0	0
5	I	1	Total 1	Zn 1	0	0
5	M	1	Total 1	Zn 1	0	0

- Molecule 6 is 2',3'-DIDEOXYCYTIDINE 5'-TRIPHOSPHATE (CCD ID: DCT) (formula: C₉H₁₆N₃O₁₂P₃).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	A	1	Total 27	C 9	N 3	O 12	P 3	0	0
6	E	1	Total 27	C 9	N 3	O 12	P 3	0	0
6	I	1	Total 27	C 9	N 3	O 12	P 3	0	0
6	M	1	Total 27	C 9	N 3	O 12	P 3	0	0

- Molecule 7 is ISOPROPYL ALCOHOL (CCD ID: IPA) (formula: $\text{C}_3\text{H}_8\text{O}$).



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

- Molecule 8 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	A	163	Total O 163 163	0	0
8	B	32	Total O 32 32	0	0
8	C	34	Total O 34 34	0	0
8	D	3	Total O 3 3	0	0
8	E	180	Total O 180 180	0	0
8	F	27	Total O 27 27	0	0
8	G	19	Total O 19 19	0	0
8	H	2	Total O 2 2	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	I	156	Total 156	O 156	0	0
8	J	29	Total 29	O 29	0	0
8	K	24	Total 24	O 24	0	0
8	L	1	Total 1	O 1	0	0
8	M	169	Total 169	O 169	0	0
8	N	29	Total 29	O 29	0	0
8	O	25	Total 25	O 25	0	0
8	P	5	Total 5	O 5	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.

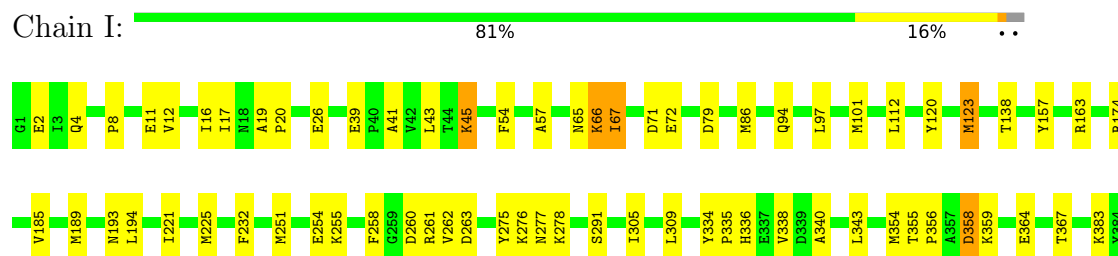
• Molecule 1: Polymerase



• Molecule 1: Polymerase



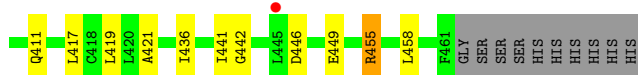
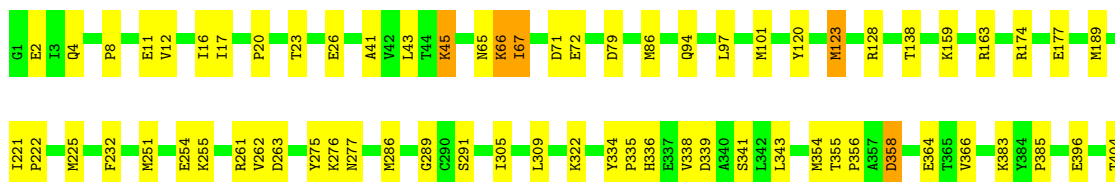
• Molecule 1: Polymerase





- Molecule 1: Polymerase

Chain M: 81% 15% ..



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

Chain B: 12% 50% 8% 31%



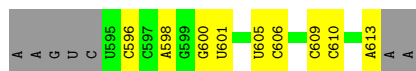
- Molecule 2: RNA (5'-R(*AP*AP*GP*UP*CP*U*CP*CP*AP*GP*GP*UP*CP*UP*CP*UP*CP*GP*UP*CP*CP*GP*GP*AP*AP*A)-3')

Chain F: 15% 46% 8% 31%



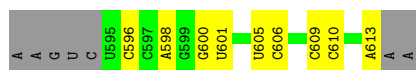
- Molecule 2: RNA (5'-R(*AP*AP*GP*UP*CP*U*CP*CP*AP*GP*GP*UP*CP*UP*CP*UP*CP*GP*UP*CP*CP*GP*GP*AP*AP*A)-3')

Chain J: 38% 35% 27%



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

Chain N: 38% 35% 27%



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

Chain C:  79% 21%



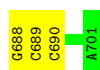
- Molecule 3: RNA (5'-R(*GP*CP*CP*CP*GP*GP*AP*CP*GP*AP*GP*AP*GP*A)-3')

Chain G:  86% 14%



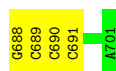
- Molecule 3: RNA (5'-R(*GP*CP*CP*CP*GP*GP*AP*CP*GP*AP*GP*AP*GP*A)-3')

Chain K:  79% 21%



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

Chain O:  71% 29%



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

Chain D:  22% 11% 67%



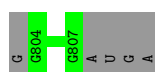
- Molecule 4: RNA (5'-R(*GP*GP*GP*AP*GP*AP*UP*GP*A)-3')

Chain H:  22% 11% 67%



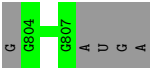
- Molecule 4: RNA (5'-R(*GP*GP*GP*AP*GP*AP*UP*GP*A)-3')

Chain L:  44% 56%



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

Chain P:  44% 56%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	60.53Å 60.54Å 192.94Å 83.82° 83.84° 77.35°	Depositor
Resolution (Å)	45.23 – 2.41 45.23 – 2.41	Depositor EDS
% Data completeness (in resolution range)	96.9 (45.23-2.41) 97.3 (45.23-2.41)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.32 (at 2.10Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.6.4_486)	Depositor
R, R_{free}	0.217 , 0.261 0.200 , 0.248	Depositor DCC
R_{free} test set	7555 reflections (4.89%)	wwPDB-VP
Wilson B-factor (Å ²)	55.0	Xtriage
Anisotropy	0.194	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 40.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.63$, $\langle L^2 \rangle = 0.50$	Xtriage
Estimated twinning fraction	0.430 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	18860	wwPDB-VP
Average B, all atoms (Å ²)	65.0	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 50.14 % 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 6.7521e-05. 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: DCT, ZN, IPA

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	0.54	0/3787	0.80	4/5122 (0.1%)
1	E	0.54	0/3787	0.81	5/5122 (0.1%)
1	I	0.55	1/3787 (0.0%)	0.80	0/5122
1	M	0.55	0/3787	0.80	1/5122 (0.0%)
2	B	0.37	0/400	0.41	0/621
2	F	0.40	0/400	0.40	0/621
2	J	0.37	0/422	0.43	0/655
2	N	0.37	0/422	0.43	0/655
3	C	0.31	0/340	0.40	0/530
3	G	0.33	0/340	0.40	0/530
3	K	0.31	0/340	0.39	0/530
3	O	0.33	0/340	0.39	0/530
4	D	0.14	0/76	0.27	0/117
4	H	0.14	0/76	0.28	0/117
4	L	0.15	0/102	0.27	0/158
4	P	0.14	0/102	0.27	0/158
All	All	0.51	1/18508 (0.0%)	0.74	10/25710 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	I	221	ILE	CA-CB	7.36	1.58	1.54

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	384	TYR	CA-C-N	5.63	125.30	119.05
1	E	384	TYR	C-N-CA	5.63	125.30	119.05
1	M	421	ALA	N-CA-C	5.58	117.45	111.36
1	E	421	ALA	N-CA-C	5.36	117.20	111.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	47	ASP	CA-C-N	5.15	125.45	119.47
1	A	47	ASP	C-N-CA	5.15	125.45	119.47
1	A	384	TYR	CA-C-N	5.08	124.69	119.05
1	A	384	TYR	C-N-CA	5.08	124.69	119.05
1	E	47	ASP	CA-C-N	5.04	125.31	119.47
1	E	47	ASP	C-N-CA	5.04	125.31	119.47

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3697	0	3658	74	0
1	E	3697	0	3658	78	0
1	I	3697	0	3658	61	0
1	M	3697	0	3658	60	0
2	B	361	0	183	20	0
2	F	361	0	183	20	0
2	J	381	0	193	5	0
2	N	381	0	193	6	0
3	C	303	0	156	5	0
3	G	303	0	156	4	0
3	K	303	0	156	5	0
3	O	303	0	156	7	0
4	D	68	0	34	0	0
4	H	68	0	34	0	0
4	L	91	0	45	0	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	27	0	12	2	0
6	E	27	0	12	3	0
6	I	27	0	12	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	M	27	0	12	3	0
7	A	8	0	16	7	0
7	E	16	0	32	6	0
7	I	12	0	24	7	0
7	M	8	0	16	9	0
7	O	4	0	8	1	0
8	A	163	0	0	5	0
8	B	32	0	0	1	0
8	C	34	0	0	0	0
8	D	3	0	0	0	0
8	E	180	0	0	7	0
8	F	27	0	0	2	0
8	G	19	0	0	0	0
8	H	2	0	0	0	0
8	I	156	0	0	8	0
8	J	29	0	0	0	0
8	K	24	0	0	0	0
8	L	1	0	0	0	0
8	M	169	0	0	7	0
8	N	29	0	0	0	0
8	O	25	0	0	1	0
8	P	5	0	0	0	0
All	All	18860	0	16310	334	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 10.

All (334) 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:364:GLU:H	7:M:6025:IPA:H13	1.05	1.07
1:E:404:THR:HA	7:E:6010:IPA:H12	1.48	0.96
1:E:70:VAL:HG11	1:E:251:MET:HE1	1.47	0.96
1:I:336:HIS:HB2	7:I:6023:IPA:H33	1.44	0.95
1:M:336:HIS:HB2	7:M:6031:IPA:C1	1.98	0.94
1:A:70:VAL:HG11	1:A:251:MET:HE1	1.47	0.94
1:M:364:GLU:N	7:M:6025:IPA:H13	1.90	0.87
1:A:336:HIS:HB2	7:A:6032:IPA:C3	2.05	0.86
3:K:688:G:H5'	2:N:613:A:OP1	1.78	0.83
1:I:258:PHE:HZ	8:I:839:HOH:O	1.61	0.82
2:F:597:C:C2'	2:F:598:A:H5'	2.09	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:597:C:C2'	2:B:598:A:H5'	2.09	0.82
2:J:613:A:OP1	3:O:688:G:H5'	1.80	0.81
1:A:411:GLN:HE22	1:A:446:ASP:H	1.25	0.81
1:E:284:GLY:HA2	8:E:1002:HOH:O	1.81	0.81
1:A:336:HIS:HB2	7:A:6032:IPA:H32	1.63	0.79
2:B:597:C:H2'	2:B:598:A:H5'	1.64	0.79
2:F:597:C:H2'	2:F:598:A:H5'	1.64	0.79
1:E:411:GLN:HE22	1:E:446:ASP:H	1.32	0.78
1:E:336:HIS:HB2	7:E:6028:IPA:C1	2.15	0.77
1:M:65:ASN:O	1:M:66:LYS:HB2	1.85	0.77
1:E:174:ARG:NH2	6:E:4003:DCT:H5	2.00	0.76
1:A:41:ALA:HB3	8:A:974:HOH:O	1.86	0.75
2:B:596:C:H3'	2:B:596:C:H6	1.51	0.75
2:F:596:C:H3'	2:F:596:C:H6	1.52	0.75
1:M:336:HIS:HB2	7:M:6031:IPA:H13	1.67	0.75
1:I:65:ASN:O	1:I:66:LYS:HB2	1.85	0.74
1:A:70:VAL:HG21	1:A:251:MET:CE	2.18	0.73
1:A:409:ASN:ND2	8:A:515:HOH:O	2.21	0.73
1:M:232:PHE:CE1	1:M:354:MET:HE3	2.24	0.73
1:I:336:HIS:HB2	7:I:6023:IPA:C3	2.18	0.72
1:E:70:VAL:HG21	1:E:251:MET:CE	2.19	0.72
1:M:336:HIS:HB2	7:M:6031:IPA:H11	1.70	0.72
3:C:688:G:H5'	2:F:613:A:O5'	1.90	0.71
1:A:70:VAL:HG21	1:A:251:MET:HE2	1.72	0.71
1:I:232:PHE:CE1	1:I:354:MET:HE3	2.26	0.71
1:A:334:TYR:CE2	7:A:6032:IPA:H31	2.26	0.71
1:A:334:TYR:HE2	7:A:6032:IPA:H31	1.55	0.70
1:E:248:ALA:HA	1:E:251:MET:HE3	1.73	0.70
1:I:411:GLN:HE22	1:I:446:ASP:H	1.39	0.70
1:E:45:LYS:HD2	1:E:54:PHE:HB3	1.73	0.70
1:E:70:VAL:HG21	1:E:251:MET:HE2	1.73	0.69
1:A:45:LYS:HD2	1:A:54:PHE:HB3	1.76	0.68
1:A:248:ALA:HA	1:A:251:MET:HE3	1.74	0.68
1:A:411:GLN:HE22	1:A:446:ASP:N	1.91	0.68
1:A:270:HIS:ND1	1:A:283:LYS:HE3	2.10	0.67
1:M:411:GLN:HE22	1:M:446:ASP:H	1.41	0.67
1:E:415:ARG:HD2	8:E:517:HOH:O	1.93	0.67
2:B:613:A:O5'	3:G:688:G:H5'	1.96	0.66
1:E:411:GLN:HE22	1:E:446:ASP:N	1.94	0.66
1:A:45:LYS:NZ	1:A:45:LYS:HB3	2.11	0.66
3:K:688:G:HO5'	3:K:688:G:H8	1.43	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:M:12:VAL:HG12	1:M:12:VAL:O	1.97	0.65
1:E:419:LEU:HD11	2:F:606:C:H4'	1.78	0.65
1:E:45:LYS:NZ	1:E:45:LYS:HB3	2.11	0.65
1:A:336:HIS:HB2	7:A:6032:IPA:H33	1.77	0.65
1:A:232:PHE:CE1	1:A:354:MET:HE3	2.33	0.64
1:I:258:PHE:CZ	8:I:839:HOH:O	2.42	0.64
3:O:688:G:H8	3:O:688:G:HO5'	1.43	0.64
1:E:270:HIS:ND1	1:E:283:LYS:HE3	2.13	0.64
1:E:232:PHE:CE1	1:E:354:MET:HE3	2.34	0.63
1:M:364:GLU:H	7:M:6025:IPA:C1	1.95	0.62
2:F:596:C:H3'	2:F:596:C:C6	2.35	0.62
1:I:12:VAL:HG12	1:I:12:VAL:O	1.99	0.62
1:M:455:ARG:HH21	1:M:458:LEU:HD12	1.63	0.62
1:A:419:LEU:HD11	2:B:606:C:H4'	1.82	0.61
1:I:411:GLN:HE22	1:I:446:ASP:N	1.98	0.61
1:M:16:ILE:HG23	1:M:277:ASN:HA	1.82	0.61
1:M:128:ARG:HG3	8:M:534:HOH:O	1.99	0.61
1:E:21:SER:HB2	1:E:44:THR:HG21	1.82	0.61
1:M:159:LYS:NZ	8:M:989:HOH:O	2.34	0.60
1:I:16:ILE:HG23	1:I:277:ASN:HA	1.83	0.60
2:B:596:C:H3'	2:B:596:C:C6	2.35	0.60
1:E:45:LYS:HD3	1:E:55:GLU:HG3	1.83	0.60
1:I:455:ARG:HH21	1:I:458:LEU:HD12	1.67	0.60
1:M:411:GLN:HE22	1:M:446:ASP:N	2.00	0.59
1:I:8:PRO:O	1:I:11:GLU:HB3	2.02	0.59
1:M:8:PRO:O	1:M:11:GLU:HB3	2.02	0.59
1:A:170:GLN:HA	1:A:170:GLN:HE21	1.66	0.59
1:A:45:LYS:HD3	1:A:55:GLU:HG3	1.84	0.59
2:B:596:C:C6	2:B:596:C:C3'	2.86	0.59
1:E:170:GLN:HE21	1:E:170:GLN:HA	1.66	0.58
1:I:193:ASN:HB2	8:I:839:HOH:O	2.03	0.58
2:F:596:C:C6	2:F:596:C:C3'	2.86	0.58
2:B:596:C:C2'	2:B:597:C:O5'	2.52	0.58
1:A:70:VAL:CG1	1:A:251:MET:HE1	2.30	0.58
1:M:383:LYS:NZ	1:M:383:LYS:HB3	2.18	0.58
1:M:336:HIS:CB	7:M:6031:IPA:H11	2.34	0.57
1:A:21:SER:HB2	1:A:44:THR:HG21	1.84	0.57
1:I:94:GLN:NE2	8:I:587:HOH:O	2.38	0.57
2:F:596:C:C2'	2:F:597:C:O5'	2.53	0.57
1:M:17:ILE:O	1:M:276:LYS:HA	2.05	0.57
1:E:294:SER:HB2	8:F:32:HOH:O	2.05	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:16:ILE:HD12	1:I:16:ILE:H	1.70	0.57
1:E:70:VAL:CG1	1:E:251:MET:HE1	2.30	0.56
1:E:139:LYS:HE2	8:E:1032:HOH:O	2.05	0.56
1:A:6:MET:SD	1:A:278:LYS:HE3	2.46	0.55
1:A:334:TYR:HE2	7:A:6032:IPA:C3	2.19	0.55
1:E:6:MET:SD	1:E:278:LYS:HE3	2.47	0.55
1:M:404:THR:HA	7:O:6009:IPA:H12	1.89	0.55
1:A:398:HIS:O	1:A:402:ARG:HG3	2.06	0.55
2:B:596:C:H2'	2:B:597:C:O5'	2.07	0.55
1:I:20:PRO:HG3	2:J:598:A:C4	2.41	0.55
1:M:436:ILE:O	1:M:442:GLY:HA3	2.07	0.55
1:E:16:ILE:HG13	1:E:277:ASN:HA	1.88	0.55
1:E:336:HIS:HB2	7:E:6028:IPA:H11	1.88	0.55
2:F:596:C:H6	2:F:596:C:C3'	2.19	0.54
1:A:294:SER:HB2	8:A:521:HOH:O	2.07	0.54
1:A:270:HIS:CE1	1:A:283:LYS:HE3	2.42	0.54
2:B:596:C:H6	2:B:596:C:C3'	2.19	0.54
1:M:16:ILE:H	1:M:16:ILE:HD12	1.72	0.54
1:M:41:ALA:HB2	1:M:163:ARG:CG	2.38	0.54
1:E:404:THR:HA	7:E:6010:IPA:C1	2.29	0.54
1:A:16:ILE:HG13	1:A:277:ASN:HA	1.90	0.54
1:E:337:GLU:O	1:E:337:GLU:HG3	2.08	0.54
1:I:41:ALA:HB2	1:I:163:ARG:CG	2.38	0.54
1:M:232:PHE:CD1	1:M:354:MET:HE3	2.43	0.54
2:F:605:U:H2'	2:F:606:C:C6	2.43	0.53
1:I:383:LYS:NZ	1:I:383:LYS:HB3	2.22	0.53
2:F:596:C:H2'	2:F:597:C:O5'	2.07	0.53
1:A:415:ARG:HD2	8:A:538:HOH:O	2.07	0.53
1:M:20:PRO:HG3	2:N:598:A:C4	2.44	0.53
1:A:20:PRO:HG3	2:B:598:A:C4	2.44	0.53
1:I:17:ILE:O	1:I:276:LYS:HA	2.09	0.53
1:M:16:ILE:HG23	1:M:277:ASN:CA	2.38	0.53
3:K:688:G:H8	3:K:688:G:O5'	1.91	0.53
1:E:71:ASP:OD1	1:E:71:ASP:C	2.52	0.53
1:I:340:ALA:HB3	7:I:6026:IPA:H33	1.89	0.53
1:E:270:HIS:CE1	1:E:283:LYS:HE3	2.44	0.52
1:I:436:ILE:O	1:I:442:GLY:HA3	2.08	0.52
2:B:609:C:O2'	2:B:610:C:H5'	2.10	0.52
1:A:4:GLN:NE2	1:A:283:LYS:HG3	2.25	0.52
1:A:71:ASP:C	1:A:71:ASP:OD1	2.52	0.52
2:F:609:C:O2'	2:F:610:C:H5'	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:232:PHE:CD1	1:I:354:MET:HE3	2.45	0.52
3:O:688:G:H8	3:O:688:G:O5'	1.92	0.52
1:M:232:PHE:CZ	1:M:354:MET:HE3	2.45	0.52
1:E:20:PRO:HG3	2:F:598:A:C4	2.44	0.51
1:A:277:ASN:OD1	1:A:278:LYS:HG3	2.10	0.51
1:A:287:PRO:O	1:A:293:THR:HG21	2.11	0.51
1:E:277:ASN:OD1	1:E:278:LYS:HG3	2.09	0.51
1:I:423:HIS:HB2	8:I:540:HOH:O	2.10	0.51
1:E:4:GLN:NE2	1:E:283:LYS:HG3	2.25	0.51
1:M:97:LEU:O	1:M:101:MET:HG3	2.11	0.51
7:M:6025:IPA:H12	8:M:604:HOH:O	2.10	0.51
1:E:398:HIS:O	1:E:402:ARG:HG3	2.10	0.51
1:I:16:ILE:HG23	1:I:277:ASN:CA	2.40	0.51
2:B:605:U:H2'	2:B:606:C:C6	2.46	0.51
1:E:16:ILE:CG1	1:E:277:ASN:HA	2.40	0.51
1:E:287:PRO:O	1:E:293:THR:HG21	2.11	0.51
2:J:600:G:O2'	2:J:601:U:H5'	2.11	0.51
1:A:2:GLU:HA	1:A:62:TYR:HB3	1.93	0.50
2:N:605:U:H2'	2:N:606:C:C6	2.45	0.50
1:A:16:ILE:CG1	1:A:277:ASN:HA	2.42	0.50
1:A:337:GLU:O	1:A:337:GLU:HG3	2.11	0.50
1:I:97:LEU:HD23	1:I:138:THR:HB	1.94	0.50
1:I:364:GLU:O	7:I:6026:IPA:H32	2.12	0.50
2:N:600:G:O2'	2:N:601:U:H5'	2.10	0.50
1:I:232:PHE:CZ	1:I:354:MET:HE3	2.47	0.50
1:E:336:HIS:HB2	7:E:6028:IPA:H13	1.92	0.49
1:M:449:GLU:OE1	8:M:553:HOH:O	2.20	0.49
1:A:41:ALA:CB	8:A:974:HOH:O	2.53	0.49
1:E:74:MET:HE1	1:E:244:ALA:HB1	1.93	0.49
1:E:404:THR:CA	7:E:6010:IPA:H12	2.33	0.49
1:A:321:LEU:O	1:A:322:LYS:HD3	2.13	0.49
1:A:270:HIS:HD1	1:A:283:LYS:HE3	1.76	0.49
1:A:334:TYR:CE2	7:A:6032:IPA:C3	2.95	0.49
1:M:383:LYS:HB3	1:M:383:LYS:HZ3	1.76	0.49
1:E:176:ILE:HD13	2:F:600:G:N7	2.28	0.49
3:O:689:C:C2'	3:O:690:C:H5'	2.43	0.49
1:A:176:ILE:HD13	2:B:600:G:N7	2.28	0.49
1:M:41:ALA:HB2	1:M:163:ARG:HG3	1.95	0.49
2:B:596:C:H2'	2:B:597:C:O4'	2.13	0.49
1:E:177:GLU:O	1:E:289:GLY:HA3	2.13	0.49
6:M:4001:DCT:H6	6:M:4001:DCT:H5'	1.95	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:M:97:LEU:HD23	1:M:138:THR:HB	1.94	0.48
2:F:607:G:H2'	2:F:608:U:C6	2.47	0.48
1:M:86:MET:HE3	1:M:86:MET:HA	1.95	0.48
6:M:4001:DCT:H5''	6:M:4001:DCT:PB	2.53	0.48
1:I:41:ALA:HB2	1:I:163:ARG:HG3	1.96	0.48
1:M:286:MET:HE2	8:M:533:HOH:O	2.13	0.48
1:I:2:GLU:HG3	1:I:4:GLN:NE2	2.28	0.48
1:A:177:GLU:O	1:A:289:GLY:HA3	2.13	0.48
2:B:607:G:H2'	2:B:608:U:C6	2.48	0.48
1:M:2:GLU:HG3	1:M:4:GLN:NE2	2.29	0.48
1:E:2:GLU:HA	1:E:62:TYR:HB3	1.95	0.48
2:F:596:C:H2'	2:F:597:C:O4'	2.13	0.48
1:I:340:ALA:H	7:I:6026:IPA:C1	2.26	0.48
6:A:4004:DCT:O4'	3:C:701:A:H2'	2.14	0.47
2:J:605:U:H2'	2:J:606:C:C6	2.49	0.47
1:A:74:MET:HE1	1:A:244:ALA:HB1	1.96	0.47
3:K:689:C:C2'	3:K:690:C:H5'	2.44	0.47
1:M:94:GLN:HG3	1:M:189:MET:O	2.14	0.47
1:A:174:ARG:NH2	6:A:4004:DCT:H5	2.30	0.47
3:C:689:C:N3	3:G:688:G:O6	2.47	0.47
1:M:309:LEU:HD23	1:M:343:LEU:HD21	1.95	0.47
6:E:4003:DCT:H3'2	8:E:656:HOH:O	2.15	0.47
1:A:70:VAL:HG21	1:A:251:MET:HE1	1.96	0.47
1:E:321:LEU:O	1:E:322:LYS:HD3	2.15	0.47
1:M:254:GLU:HG3	1:M:262:VAL:HG11	1.97	0.47
1:A:237:TYR:CG	1:A:328:ASP:HB3	2.50	0.47
2:B:600:G:O2'	2:B:601:U:H5'	2.15	0.47
2:F:600:G:O2'	2:F:601:U:H5'	2.15	0.47
3:C:688:G:O6	3:G:689:C:N3	2.48	0.46
1:E:339:ASP:OD1	1:E:341:SER:OG	2.26	0.46
1:A:165:LYS:HB3	1:A:165:LYS:HE2	1.62	0.46
1:I:309:LEU:HD23	1:I:343:LEU:HD21	1.97	0.46
1:E:159:LYS:HD2	1:E:176:ILE:HD11	1.97	0.46
1:I:355:THR:HB	1:I:356:PRO:HD2	1.98	0.46
1:A:440:PRO:HA	1:A:443:ARG:NH1	2.30	0.46
1:A:79:ASP:OD1	1:A:255:LYS:HE2	2.16	0.46
1:I:79:ASP:OD1	1:I:255:LYS:HE2	2.16	0.46
2:B:599:G:C8	8:B:1028:HOH:O	2.56	0.46
1:M:41:ALA:HB2	1:M:163:ARG:HG2	1.98	0.46
1:A:304:ILE:O	1:A:308:LEU:HG	2.15	0.46
1:E:174:ARG:CZ	6:E:4003:DCT:H5	2.44	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:G:689:C:H6	3:G:689:C:O5'	1.99	0.46
1:M:12:VAL:O	1:M:12:VAL:CG1	2.63	0.46
1:A:440:PRO:HA	1:A:443:ARG:HH11	1.80	0.46
1:I:41:ALA:HB2	1:I:163:ARG:HG2	1.97	0.46
1:I:86:MET:HE3	1:I:86:MET:HA	1.98	0.46
1:I:94:GLN:HG3	1:I:189:MET:O	2.16	0.46
1:E:419:LEU:HD11	2:F:606:C:C4'	2.46	0.45
1:A:8:PRO:HD2	1:A:11:GLU:OE2	2.17	0.45
1:A:105:ASP:OD2	1:A:200:LYS:HE2	2.16	0.45
1:E:237:TYR:CG	1:E:328:ASP:HB3	2.51	0.45
1:I:334:TYR:CD2	1:I:335:PRO:HD2	2.51	0.45
1:M:120:TYR:HA	1:M:123:MET:HG3	1.97	0.45
1:E:237:TYR:CD2	1:E:328:ASP:HB3	2.52	0.45
1:E:105:ASP:OD2	1:E:200:LYS:HE2	2.17	0.45
1:M:97:LEU:CD2	1:M:138:THR:HB	2.46	0.45
3:C:689:C:O5'	3:C:689:C:H6	1.99	0.45
1:I:12:VAL:O	1:I:12:VAL:CG1	2.64	0.45
1:I:358:ASP:OD1	1:I:358:ASP:N	2.49	0.45
1:M:334:TYR:CD2	1:M:335:PRO:HD2	2.52	0.45
3:O:689:C:O5'	3:O:689:C:H6	2.00	0.45
1:A:431:LYS:HB2	1:A:431:LYS:HE2	1.74	0.45
1:E:108:GLU:HG2	8:F:265:HOH:O	2.16	0.45
1:M:45:LYS:NZ	1:M:275:TYR:OH	2.49	0.45
1:E:309:LEU:HD23	1:E:343:LEU:HD21	1.98	0.45
1:I:120:TYR:HA	1:I:123:MET:HG3	1.99	0.45
1:A:5:TRP:O	1:A:280:TYR:HA	2.16	0.44
1:E:5:TRP:O	1:E:280:TYR:HA	2.17	0.44
1:E:232:PHE:CD1	1:E:354:MET:HE3	2.52	0.44
1:M:71:ASP:C	1:M:71:ASP:OD1	2.60	0.44
1:A:237:TYR:CD2	1:A:328:ASP:HB3	2.52	0.44
1:I:254:GLU:HG3	1:I:262:VAL:HG11	2.00	0.44
1:A:447:LEU:HA	1:A:448:PRO:HD3	1.81	0.44
1:M:358:ASP:N	1:M:358:ASP:OD1	2.50	0.44
1:M:417:LEU:HD23	1:M:417:LEU:HA	1.84	0.43
1:I:66:LYS:HB3	1:I:67:ILE:H	1.55	0.43
6:M:4001:DCT:H5''	6:M:4001:DCT:O1B	2.18	0.43
1:E:440:PRO:HA	1:E:443:ARG:HH11	1.82	0.43
1:I:336:HIS:CB	7:I:6023:IPA:C3	2.94	0.43
3:K:689:C:H6	3:K:689:C:O5'	2.01	0.43
1:M:177:GLU:O	1:M:289:GLY:HA3	2.19	0.43
3:O:689:C:OP2	8:O:730:HOH:O	2.21	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:70:VAL:HG21	1:E:251:MET:HE1	1.99	0.43
1:E:269:ASN:OD1	8:E:1002:HOH:O	2.21	0.43
1:A:406:ASP:HB3	1:A:409:ASN:ND2	2.33	0.43
1:E:97:LEU:O	1:E:101:MET:HG3	2.18	0.43
1:I:97:LEU:O	1:I:101:MET:HG3	2.18	0.43
1:E:8:PRO:HD2	1:E:11:GLU:OE2	2.18	0.43
1:E:37:VAL:HG22	1:E:399:GLU:OE2	2.18	0.43
1:A:232:PHE:CD1	1:A:354:MET:HE3	2.54	0.43
1:A:309:LEU:HD23	1:A:343:LEU:HD21	2.00	0.43
1:M:322:LYS:HA	1:M:322:LYS:HD3	1.82	0.43
8:M:520:HOH:O	2:N:600:G:H4'	2.18	0.43
1:I:45:LYS:NZ	1:I:275:TYR:OH	2.51	0.43
1:E:18:ASN:HD22	1:E:18:ASN:HA	1.62	0.43
1:E:170:GLN:HA	1:E:170:GLN:NE2	2.33	0.43
1:E:440:PRO:HA	1:E:443:ARG:NH1	2.33	0.43
1:I:367:THR:HA	7:I:6015:IPA:H11	1.99	0.43
1:E:447:LEU:HA	1:E:448:PRO:HD3	1.81	0.43
1:M:355:THR:HB	1:M:356:PRO:HD2	2.01	0.43
1:E:139:LYS:HE3	1:E:139:LYS:HB2	1.87	0.42
1:A:170:GLN:HA	1:A:170:GLN:NE2	2.33	0.42
1:I:71:ASP:OD1	1:I:71:ASP:C	2.61	0.42
1:A:67:ILE:HG13	1:A:244:ALA:HB3	2.02	0.42
2:B:602:C:H2'	2:B:603:U:C6	2.55	0.42
1:I:193:ASN:CB	8:I:839:HOH:O	2.63	0.42
1:I:359:LYS:HA	8:I:514:HOH:O	2.19	0.42
1:I:97:LEU:CD2	1:I:138:THR:HB	2.49	0.42
1:I:163:ARG:NH2	6:I:4002:DCT:O1G	2.41	0.42
1:I:225:MET:HE2	1:I:225:MET:HA	2.01	0.42
1:E:22:LYS:HA	8:E:1047:HOH:O	2.20	0.42
1:E:232:PHE:CZ	1:E:354:MET:HE3	2.55	0.42
1:E:270:HIS:HD1	1:E:283:LYS:HE3	1.81	0.42
1:E:345:GLN:HB2	8:E:540:HOH:O	2.20	0.42
1:A:286:MET:HE2	1:A:286:MET:HB3	1.75	0.42
1:A:232:PHE:CZ	1:A:354:MET:HE3	2.54	0.42
1:A:419:LEU:HD11	2:B:606:C:C4'	2.49	0.42
1:I:185:VAL:HG12	1:I:189:MET:HE2	2.02	0.42
2:J:609:C:O2'	2:J:610:C:H5'	2.20	0.42
1:E:304:ILE:O	1:E:308:LEU:HG	2.19	0.41
1:E:185:VAL:HG12	1:E:189:MET:HE2	2.02	0.41
1:E:7:ARG:O	1:E:8:PRO:C	2.63	0.41
1:M:66:LYS:HB3	1:M:67:ILE:H	1.53	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:37:VAL:HG22	1:A:399:GLU:OE2	2.20	0.41
2:F:602:C:H2'	2:F:603:U:C6	2.55	0.41
1:I:19:ALA:HB2	1:I:157:TYR:CD1	2.56	0.41
1:M:339:ASP:OD1	1:M:341:SER:OG	2.28	0.41
1:M:251:MET:O	1:M:255:LYS:HG3	2.19	0.41
3:O:690:C:H2'	3:O:691:C:O4'	2.21	0.41
1:A:18:ASN:HD22	1:A:18:ASN:HA	1.63	0.41
1:A:381:ASP:HB3	1:A:384:TYR:O	2.19	0.41
1:E:17:ILE:O	1:E:276:LYS:HA	2.21	0.41
1:E:381:ASP:HB3	1:E:384:TYR:O	2.21	0.41
1:A:57:ALA:O	1:A:60:SER:HB3	2.20	0.41
2:F:607:G:H2'	2:F:608:U:H6	1.86	0.41
1:I:260:ASP:OD2	1:I:260:ASP:N	2.54	0.41
1:A:7:ARG:O	1:A:8:PRO:C	2.62	0.41
1:E:54:PHE:O	1:E:57:ALA:N	2.53	0.41
1:M:79:ASP:OD1	1:M:255:LYS:HE2	2.21	0.41
2:N:609:C:O2'	2:N:610:C:H5'	2.20	0.41
1:A:154:LEU:HD21	1:A:290:CYS:SG	2.61	0.41
1:E:94:GLN:HE21	1:E:94:GLN:HB3	1.68	0.41
1:E:254:GLU:HG3	1:E:262:VAL:HG11	2.03	0.41
1:I:112:LEU:HA	1:I:112:LEU:HD23	1.80	0.41
1:I:251:MET:O	1:I:255:LYS:HG3	2.20	0.41
1:I:277:ASN:OD1	1:I:278:LYS:HG2	2.21	0.41
1:I:383:LYS:O	1:I:385:PRO:HD3	2.21	0.41
1:M:23:THR:N	8:M:801:HOH:O	2.54	0.41
1:I:54:PHE:O	1:I:57:ALA:N	2.55	0.41
1:A:339:ASP:OD1	1:A:341:SER:OG	2.30	0.40
1:M:225:MET:HE2	1:M:225:MET:HA	2.01	0.40
1:M:366:VAL:HG22	7:M:6025:IPA:H33	2.03	0.40
1:A:159:LYS:HD2	1:A:176:ILE:HD11	2.02	0.40
1:E:436:ILE:O	1:E:442:GLY:HA3	2.21	0.40
1:I:194:LEU:N	8:I:839:HOH:O	2.54	0.40
1:M:45:LYS:N	1:M:45:LYS:HD3	2.36	0.40
1:M:383:LYS:O	1:M:385:PRO:HD3	2.21	0.40
1:E:6:MET:O	1:E:7:ARG:HB3	2.22	0.40
1:M:221:ILE:HB	1:M:222:PRO:HD3	2.04	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%)	443 (96%)	15 (3%)	1 (0%)	43	57
1	E	459/471 (98%)	445 (97%)	13 (3%)	1 (0%)	43	57
1	I	459/471 (98%)	447 (97%)	11 (2%)	1 (0%)	43	57
1	M	459/471 (98%)	446 (97%)	12 (3%)	1 (0%)	43	57
All	All	1836/1884 (98%)	1781 (97%)	51 (3%)	4 (0%)	43	57

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	I	66	LYS
1	M	66	LYS
1	A	285	GLY
1	E	285	GLY

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%)	382 (95%)	21 (5%)	21	34
1	E	403/412 (98%)	381 (94%)	22 (6%)	19	32
1	I	403/412 (98%)	385 (96%)	18 (4%)	24	40
1	M	403/412 (98%)	386 (96%)	17 (4%)	26	43
All	All	1612/1648 (98%)	1534 (95%)	78 (5%)	23	38

All (78) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	GLN
1	A	16	ILE
1	A	18	ASN
1	A	22	LYS
1	A	43	LEU
1	A	45	LYS
1	A	46	ASN
1	A	60	SER
1	A	89	ASP
1	A	94	GLN
1	A	125	LYS
1	A	166	THR
1	A	170	GLN
1	A	261	ARG
1	A	269	ASN
1	A	294	SER
1	A	337	GLU
1	A	338	VAL
1	A	360	SER
1	A	396	GLU
1	A	411	GLN
1	E	4	GLN
1	E	16	ILE
1	E	18	ASN
1	E	22	LYS
1	E	43	LEU
1	E	45	LYS
1	E	46	ASN
1	E	60	SER
1	E	89	ASP
1	E	94	GLN
1	E	125	LYS
1	E	166	THR
1	E	170	GLN
1	E	261	ARG
1	E	269	ASN
1	E	291	SER
1	E	294	SER
1	E	337	GLU
1	E	338	VAL
1	E	360	SER
1	E	396	GLU

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Mol	Chain	Res	Type
1	E	411	GLN
1	I	26	GLU
1	I	39	GLU
1	I	43	LEU
1	I	45	LYS
1	I	67	ILE
1	I	72	GLU
1	I	123	MET
1	I	174	ARG
1	I	261	ARG
1	I	263	ASP
1	I	291	SER
1	I	305	ILE
1	I	338	VAL
1	I	358	ASP
1	I	396	GLU
1	I	419	LEU
1	I	441	ILE
1	I	455	ARG
1	M	26	GLU
1	M	43	LEU
1	M	45	LYS
1	M	67	ILE
1	M	72	GLU
1	M	123	MET
1	M	174	ARG
1	M	261	ARG
1	M	263	ASP
1	M	291	SER
1	M	305	ILE
1	M	338	VAL
1	M	358	ASP
1	M	396	GLU
1	M	419	LEU
1	M	441	ILE
1	M	455	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) 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	320	HIS
1	A	336	HIS
1	A	409	ASN
1	A	411	GLN
1	E	18	ASN
1	E	94	GLN
1	E	170	GLN
1	E	320	HIS
1	E	409	ASN
1	E	411	GLN
1	I	4	GLN
1	I	142	GLN
1	I	269	ASN
1	I	409	ASN
1	I	411	GLN
1	M	4	GLN
1	M	269	ASN
1	M	409	ASN
1	M	411	GLN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	B	16/26 (61%)	2 (12%)	0
2	F	16/26 (61%)	2 (12%)	0
2	J	17/26 (65%)	1 (5%)	0
2	N	17/26 (65%)	1 (5%)	0
3	C	13/14 (92%)	0	0
3	G	13/14 (92%)	0	0
3	K	13/14 (92%)	0	0
3	O	13/14 (92%)	0	0
4	D	2/9 (22%)	1 (50%)	0
4	H	2/9 (22%)	1 (50%)	0
4	L	3/9 (33%)	0	0
4	P	3/9 (33%)	0	0
All	All	128/196 (65%)	8 (6%)	0

All (8) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
2	B	597	C
2	B	598	A
4	D	806	A
2	F	597	C
2	F	598	A
4	H	806	A
2	J	596	C
2	N	596	C

There are no RNA pucker outliers to report.

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 20 ligands modelled in this entry, 4 are monoatomic - leaving 16 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	$\# Z > 2$
6	DCT	M	4001	-	27,28,28	1.84	3 (11%)	37,43,43	1.14	4 (10%)
7	IPA	E	6028	-	3,3,3	0.55	0	3,3,3	0.33	0
7	IPA	M	6025	-	3,3,3	0.54	0	3,3,3	0.25	0
6	DCT	A	4004	-	27,28,28	1.84	4 (14%)	37,43,43	1.30	4 (10%)
7	IPA	A	6008	-	3,3,3	0.61	0	3,3,3	0.35	0
7	IPA	I	6026	-	3,3,3	0.63	0	3,3,3	0.29	0
7	IPA	I	6023	-	3,3,3	0.49	0	3,3,3	0.43	0
6	DCT	E	4003	-	27,28,28	1.81	4 (14%)	37,43,43	1.22	2 (5%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	DCT	I	4002	-	27,28,28	1.82	4 (14%)	37,43,43	1.34	5 (13%)
7	IPA	A	6032	-	3,3,3	0.38	0	3,3,3	0.56	0
7	IPA	E	6010	-	3,3,3	0.47	0	3,3,3	0.42	0
7	IPA	E	6027	-	3,3,3	0.52	0	3,3,3	0.34	0
7	IPA	I	6015	-	3,3,3	0.54	0	3,3,3	0.36	0
7	IPA	E	6007	-	3,3,3	0.62	0	3,3,3	0.36	0
7	IPA	M	6031	-	3,3,3	0.29	0	3,3,3	0.61	0
7	IPA	O	6009	-	3,3,3	0.58	0	3,3,3	0.42	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
6	DCT	A	4004	-	-	3/22/31/31	0/2/2/2
6	DCT	I	4002	-	-	7/22/31/31	0/2/2/2
6	DCT	M	4001	-	-	2/22/31/31	0/2/2/2
6	DCT	E	4003	-	-	5/22/31/31	0/2/2/2

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	I	4002	DCT	O2-C2	7.11	1.37	1.23
6	M	4001	DCT	O2-C2	7.02	1.36	1.23
6	A	4004	DCT	O2-C2	6.90	1.36	1.23
6	E	4003	DCT	O2-C2	6.76	1.36	1.23
6	M	4001	DCT	PB-O3A	2.90	1.62	1.59
6	E	4003	DCT	C2-N1	-2.79	1.34	1.40
6	A	4004	DCT	C2-N1	-2.76	1.34	1.40
6	M	4001	DCT	C2-N1	-2.71	1.34	1.40
6	I	4002	DCT	C2-N1	-2.46	1.34	1.40
6	A	4004	DCT	PB-O3A	2.31	1.62	1.59
6	E	4003	DCT	PA-O3A	2.16	1.61	1.59
6	A	4004	DCT	PB-O3B	2.13	1.61	1.59
6	E	4003	DCT	PB-O3A	2.10	1.61	1.59
6	I	4002	DCT	PA-O3A	2.08	1.61	1.59
6	I	4002	DCT	PB-O3A	2.07	1.61	1.59

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	I	4002	DCT	C2'-C1'-N1	-4.11	104.61	112.40
6	A	4004	DCT	C2'-C1'-N1	-3.58	105.61	112.40
6	A	4004	DCT	C3'-C2'-C1'	3.04	106.38	102.87
6	I	4002	DCT	C3'-C2'-C1'	2.95	106.27	102.87
6	E	4003	DCT	C3'-C2'-C1'	2.81	106.11	102.87
6	I	4002	DCT	O4'-C1'-N1	2.77	112.79	107.86
6	M	4001	DCT	C3'-C2'-C1'	2.66	105.94	102.87
6	M	4001	DCT	O2B-PB-O3A	2.56	114.19	107.27
6	A	4004	DCT	O4'-C1'-N1	2.41	112.15	107.86
6	A	4004	DCT	C5-C6-N1	-2.29	118.11	121.84
6	I	4002	DCT	C5-C6-N1	-2.28	118.13	121.84
6	M	4001	DCT	C2'-C1'-N1	-2.20	108.24	112.40
6	E	4003	DCT	O2B-PB-O3A	2.10	112.95	107.27
6	M	4001	DCT	O2G-PG-O3B	2.07	111.58	104.64
6	I	4002	DCT	O4'-C4'-C5'	2.02	112.98	109.34

There are no chirality outliers.

All (17) torsion outliers are listed below:

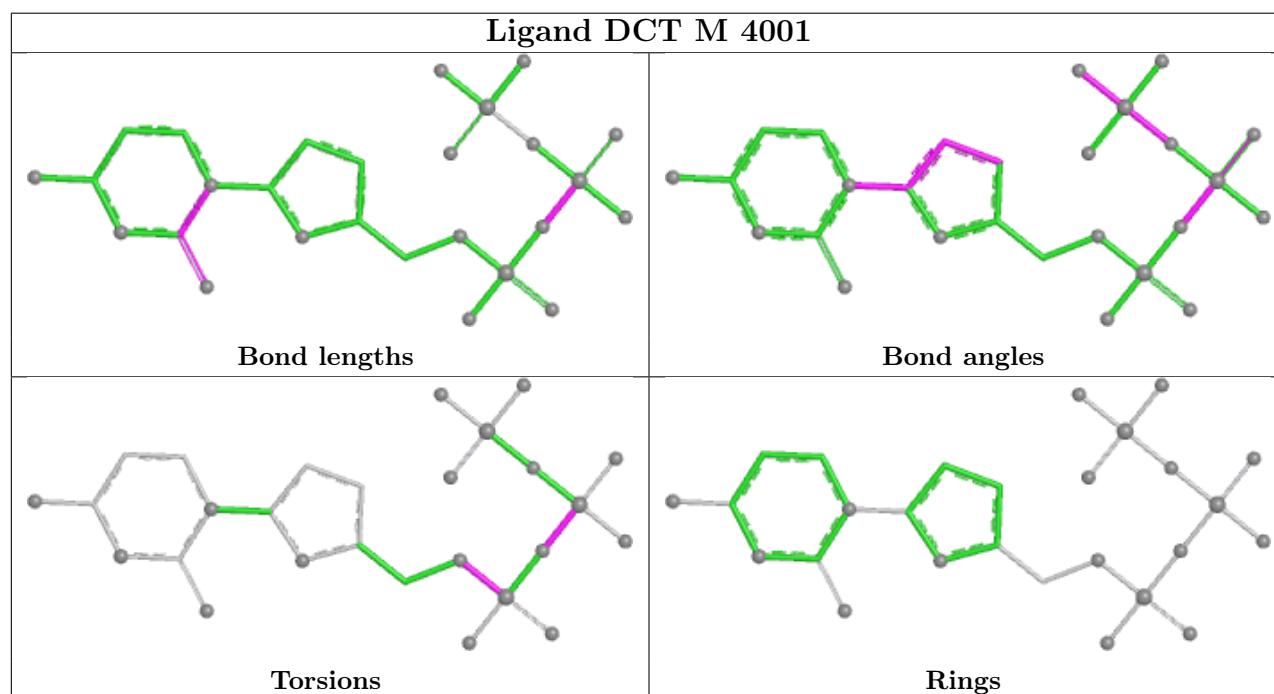
Mol	Chain	Res	Type	Atoms
6	A	4004	DCT	C5'-O5'-PA-O1A
6	A	4004	DCT	C5'-O5'-PA-O2A
6	A	4004	DCT	C5'-O5'-PA-O3A
6	E	4003	DCT	C5'-O5'-PA-O1A
6	E	4003	DCT	C5'-O5'-PA-O3A
6	I	4002	DCT	C3'-C4'-C5'-O5'
6	M	4001	DCT	C5'-O5'-PA-O1A
6	I	4002	DCT	O4'-C4'-C5'-O5'
6	I	4002	DCT	PB-O3A-PA-O1A
6	E	4003	DCT	PA-O3A-PB-O2B
6	I	4002	DCT	PG-O3B-PB-O2B
6	I	4002	DCT	PA-O3A-PB-O2B
6	E	4003	DCT	PB-O3A-PA-O2A
6	E	4003	DCT	PA-O3A-PB-O1B
6	I	4002	DCT	PB-O3A-PA-O2A
6	I	4002	DCT	PG-O3B-PB-O1B
6	M	4001	DCT	PA-O3A-PB-O1B

There are no ring outliers.

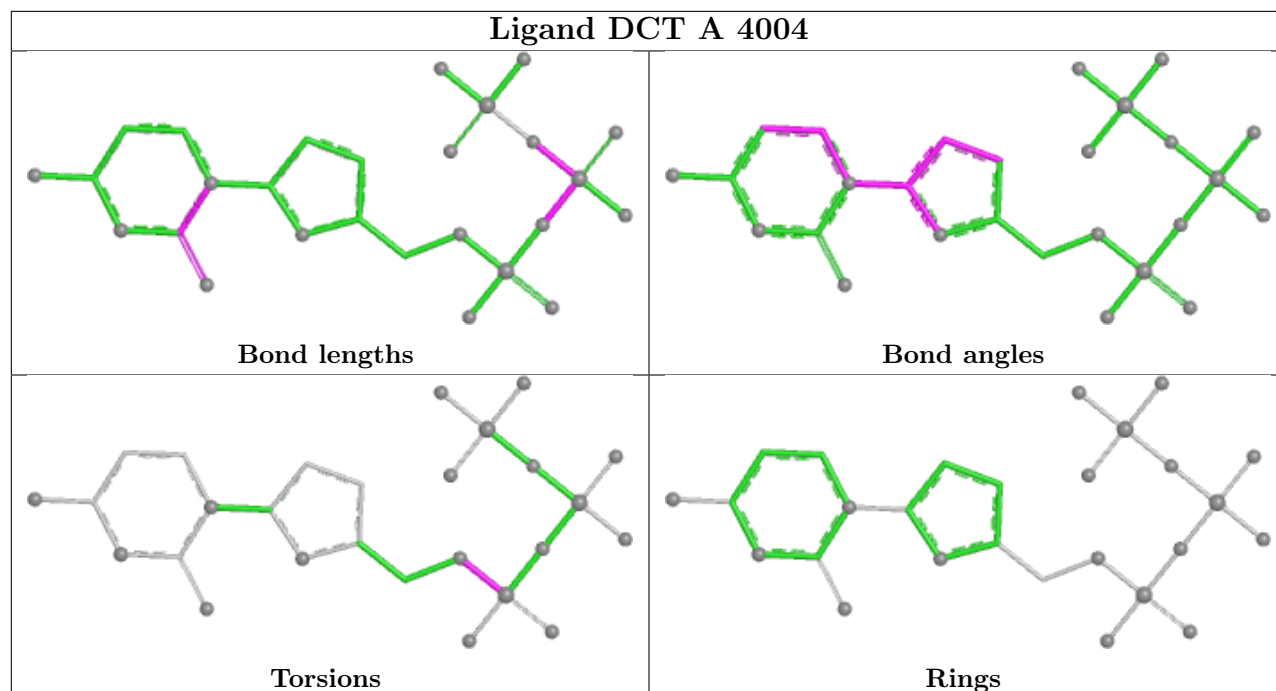
13 monomers are involved in 39 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	M	4001	DCT	3	0
7	E	6028	IPA	3	0
7	M	6025	IPA	5	0
6	A	4004	DCT	2	0
7	I	6026	IPA	3	0
7	I	6023	IPA	3	0
6	E	4003	DCT	3	0
6	I	4002	DCT	1	0
7	A	6032	IPA	7	0
7	E	6010	IPA	3	0
7	I	6015	IPA	1	0
7	M	6031	IPA	4	0
7	O	6009	IPA	1	0

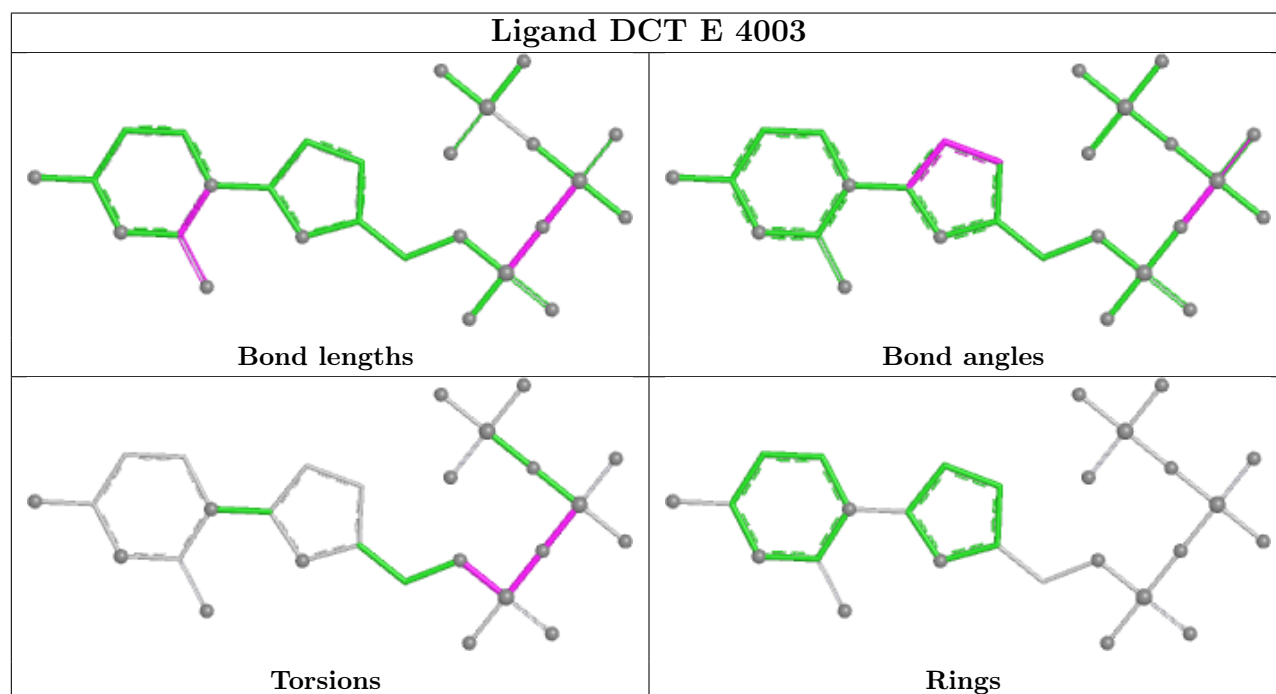
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

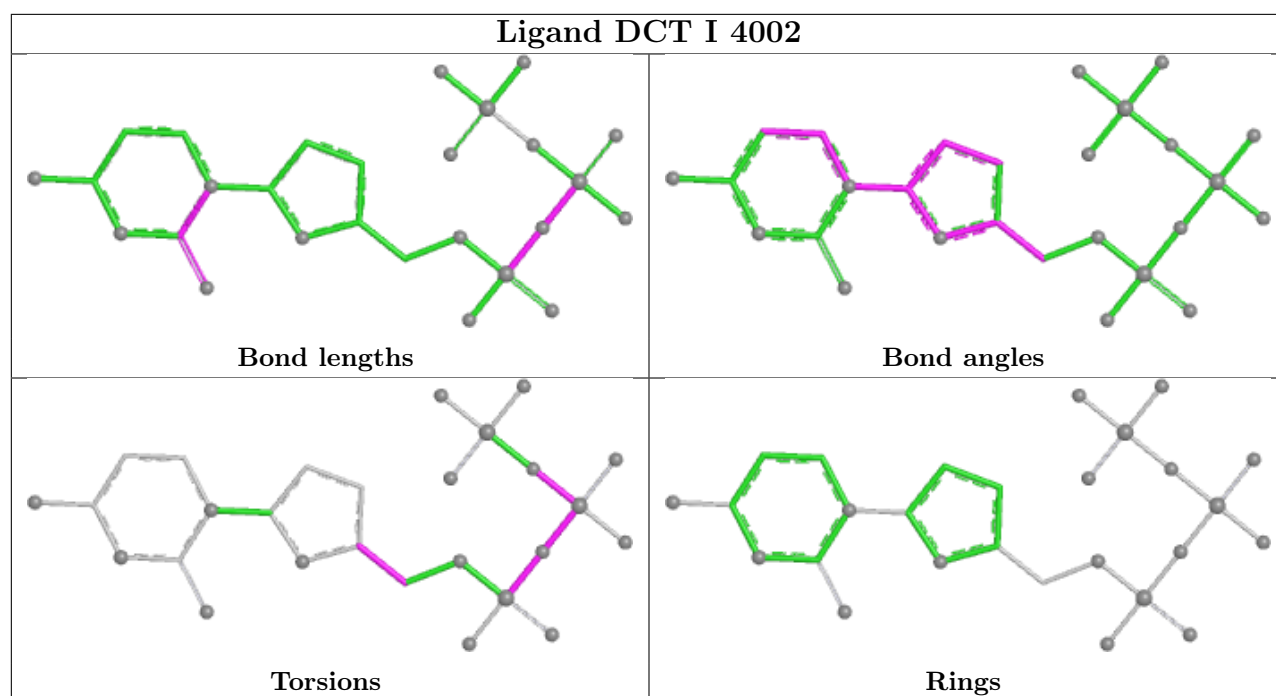


Ligand DCT A 4004



Ligand DCT E 4003





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	461/471 (97%)	-1.06	0 100 100	38, 57, 99, 130	0
1	E	461/471 (97%)	-1.08	0 100 100	37, 56, 99, 130	0
1	I	461/471 (97%)	-1.11	0 100 100	37, 57, 94, 131	0
1	M	461/471 (97%)	-1.08	1 (0%) 91 90	37, 57, 93, 131	0
2	B	18/26 (69%)	-1.62	0 100 100	47, 58, 140, 153	0
2	F	18/26 (69%)	-1.68	0 100 100	46, 57, 140, 153	0
2	J	19/26 (73%)	-1.48	0 100 100	44, 62, 158, 162	0
2	N	19/26 (73%)	-1.61	0 100 100	45, 63, 158, 162	0
3	C	14/14 (100%)	-1.78	0 100 100	47, 55, 112, 119	0
3	G	14/14 (100%)	-1.66	0 100 100	47, 55, 111, 119	0
3	K	14/14 (100%)	-1.65	0 100 100	44, 58, 120, 121	0
3	O	14/14 (100%)	-1.58	0 100 100	44, 58, 120, 121	0
4	D	3/9 (33%)	-0.84	0 100 100	134, 134, 137, 153	0
4	H	3/9 (33%)	-1.01	0 100 100	134, 134, 138, 154	0
4	L	4/9 (44%)	-1.37	0 100 100	118, 121, 137, 168	0
4	P	4/9 (44%)	-1.28	0 100 100	118, 121, 137, 168	0
All	All	1988/2080 (95%)	-1.12	1 (0%) 92 91	37, 57, 102, 168	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	M	445	LEU	2.2

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 ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

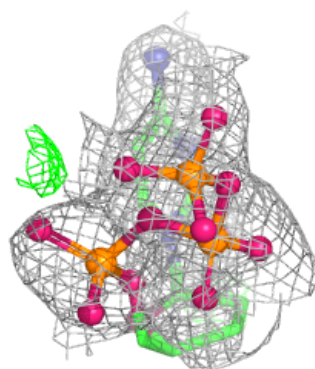
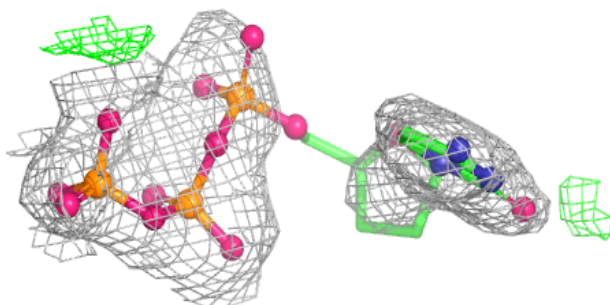
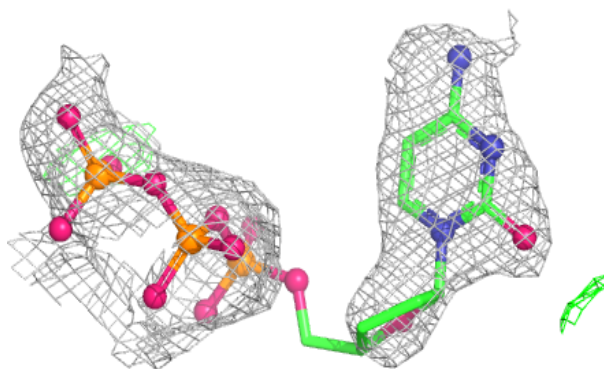
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
7	IPA	E	6027	4/4	0.87	0.18	80,81,87,88	0
7	IPA	I	6015	4/4	0.96	0.16	63,80,80,88	0
7	IPA	I	6026	4/4	0.96	0.10	49,72,82,97	0
7	IPA	E	6007	4/4	0.97	0.11	52,61,71,72	0
7	IPA	A	6032	4/4	0.97	0.12	57,78,78,86	0
5	ZN	E	2002	1/1	0.98	0.06	116,116,116,116	1
6	DCT	A	4004	27/27	0.98	0.06	35,86,154,172	27
7	IPA	A	6008	4/4	0.98	0.10	48,67,69,80	0
5	ZN	A	2001	1/1	0.98	0.07	102,102,102,102	1
7	IPA	M	6025	4/4	0.98	0.07	76,79,96,118	0
7	IPA	O	6009	4/4	0.98	0.11	55,78,78,82	0
7	IPA	E	6010	4/4	0.99	0.07	54,62,62,89	0
6	DCT	I	4002	27/27	0.99	0.05	34,74,116,186	27
7	IPA	E	6028	4/4	0.99	0.08	67,76,82,84	0
6	DCT	M	4001	27/27	0.99	0.04	36,78,139,174	27
7	IPA	I	6023	4/4	0.99	0.07	68,97,99,114	0
5	ZN	M	2004	1/1	0.99	0.02	93,93,93,93	1
5	ZN	I	2003	1/1	0.99	0.03	96,96,96,96	1
6	DCT	E	4003	27/27	0.99	0.04	36,86,138,159	27
7	IPA	M	6031	4/4	1.00	0.06	94,94,102,114	0

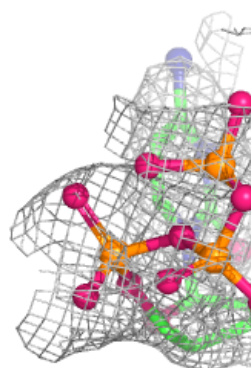
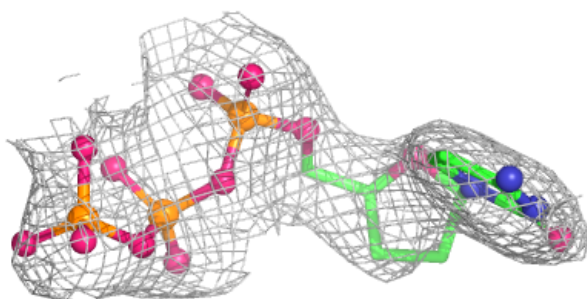
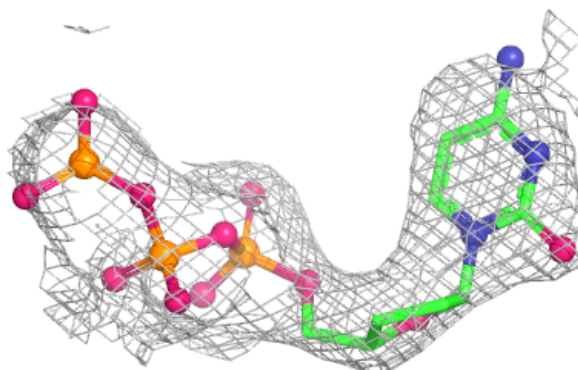
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around DCT A 4004:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

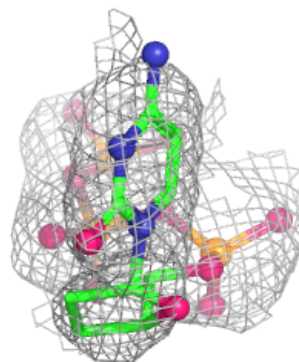
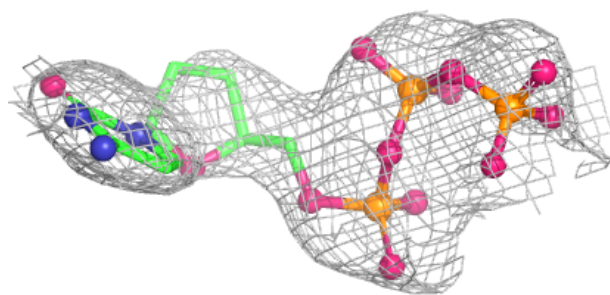
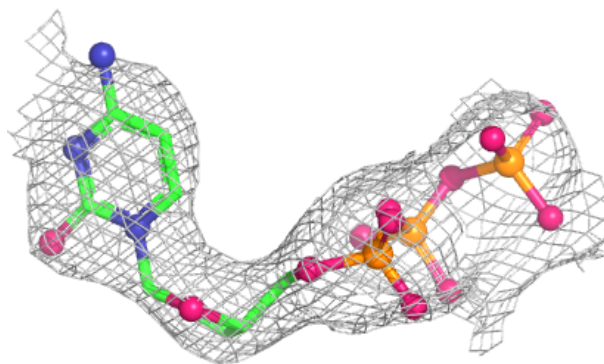
**Electron density around DCT I 4002:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

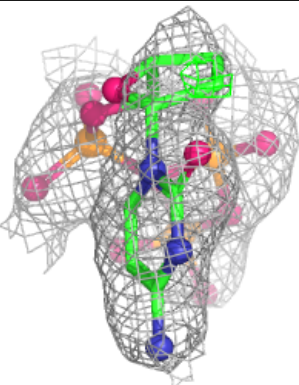
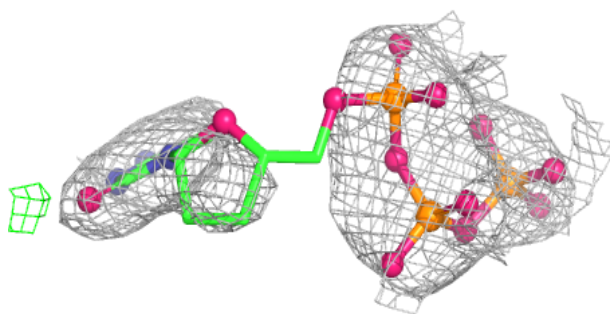
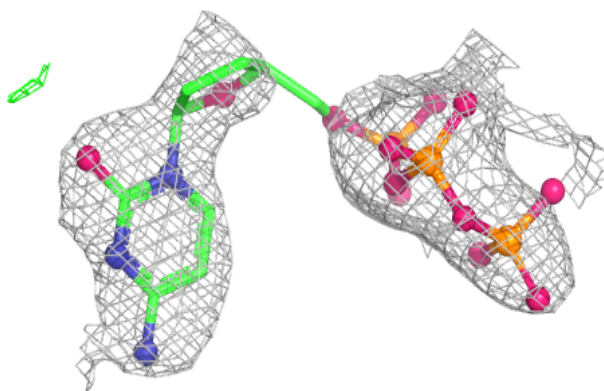


Electron density around DCT M 4001:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

**Electron density around DCT E 4003:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.