



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 10:24 AM UTC

PDB ID : 4K4G / pdb_00004k4g
Title : Ternary crystal structures of human DNA POLYMERASE LAMBDA IN COMPLEX WITH DNA AND L-DCTP.
Authors : Vyas, R.; Suo, Z.
Deposited on : 2013-04-12
Resolution : 2.15 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

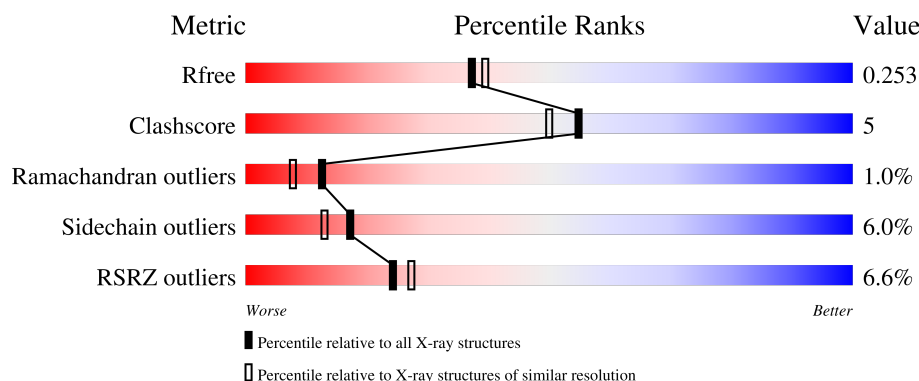
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.15 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.



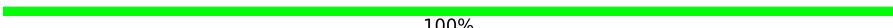
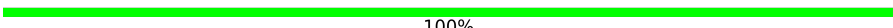
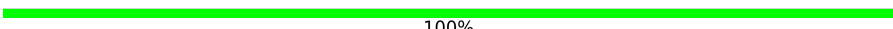
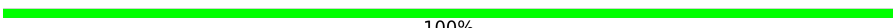
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	2057 (2.16-2.16)
Clashscore	190562	2159 (2.16-2.16)
Ramachandran outliers	187476	2134 (2.16-2.16)
Sidechain outliers	187428	2133 (2.16-2.16)
RSRZ outliers	180081	2059 (2.16-2.16)

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	340	6% 78% 15% • 5%
1	E	340	7% 77% 15% • 5%
1	I	340	6% 86% 9% • •
1	M	340	8% 80% 12% • • •
2	B	11	64% 36%

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Mol	Chain	Length	Quality of chain
2	F	11	 45%55%
2	J	11	 82%18%
2	N	11	 64%36%
3	C	6	 83%17%
3	G	6	 83%17%
3	K	6	 83%17%
3	O	6	 100%
4	D	4	 50%50%
4	H	4	 100%
4	L	4	 100%
4	P	4	 100%

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	CAC	M	607	-	X	-	-

2 Entry composition

There are 9 unique types of molecules in this entry. The entry contains 12538 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called DNA polymerase lambda.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	322	Total	C	N	O	S	0	0	0
			2548	1603	466	468	11			
1	E	323	Total	C	N	O	S	0	0	0
			2541	1600	466	463	12			
1	I	327	Total	C	N	O	S	0	0	0
			2580	1621	474	473	12			
1	M	328	Total	C	N	O	S	0	0	0
			2584	1624	474	474	12			

There are 36 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	244	MET	-	expression tag	UNP Q9UGP5
A	576	LEU	-	expression tag	UNP Q9UGP5
A	577	GLU	-	expression tag	UNP Q9UGP5
A	578	HIS	-	expression tag	UNP Q9UGP5
A	579	HIS	-	expression tag	UNP Q9UGP5
A	580	HIS	-	expression tag	UNP Q9UGP5
A	581	HIS	-	expression tag	UNP Q9UGP5
A	582	HIS	-	expression tag	UNP Q9UGP5
A	583	HIS	-	expression tag	UNP Q9UGP5
E	244	MET	-	expression tag	UNP Q9UGP5
E	576	LEU	-	expression tag	UNP Q9UGP5
E	577	GLU	-	expression tag	UNP Q9UGP5
E	578	HIS	-	expression tag	UNP Q9UGP5
E	579	HIS	-	expression tag	UNP Q9UGP5
E	580	HIS	-	expression tag	UNP Q9UGP5
E	581	HIS	-	expression tag	UNP Q9UGP5
E	582	HIS	-	expression tag	UNP Q9UGP5
E	583	HIS	-	expression tag	UNP Q9UGP5
I	244	MET	-	expression tag	UNP Q9UGP5
I	576	LEU	-	expression tag	UNP Q9UGP5
I	577	GLU	-	expression tag	UNP Q9UGP5

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Chain	Residue	Modelled	Actual	Comment	Reference
I	578	HIS	-	expression tag	UNP Q9UGP5
I	579	HIS	-	expression tag	UNP Q9UGP5
I	580	HIS	-	expression tag	UNP Q9UGP5
I	581	HIS	-	expression tag	UNP Q9UGP5
I	582	HIS	-	expression tag	UNP Q9UGP5
I	583	HIS	-	expression tag	UNP Q9UGP5
M	244	MET	-	expression tag	UNP Q9UGP5
M	576	LEU	-	expression tag	UNP Q9UGP5
M	577	GLU	-	expression tag	UNP Q9UGP5
M	578	HIS	-	expression tag	UNP Q9UGP5
M	579	HIS	-	expression tag	UNP Q9UGP5
M	580	HIS	-	expression tag	UNP Q9UGP5
M	581	HIS	-	expression tag	UNP Q9UGP5
M	582	HIS	-	expression tag	UNP Q9UGP5
M	583	HIS	-	expression tag	UNP Q9UGP5

- Molecule 2 is a DNA chain called DNA (5'-D(*CP*GP*GP*CP*GP*GP*TP*AP*CP*TP*G)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	11	Total	C	N	O	P	0	0	0
			225	107	43	65	10			
2	F	11	Total	C	N	O	P	0	0	0
			225	107	43	65	10			
2	J	11	Total	C	N	O	P	0	0	0
			225	107	43	65	10			
2	N	11	Total	C	N	O	P	0	0	0
			225	107	43	65	10			

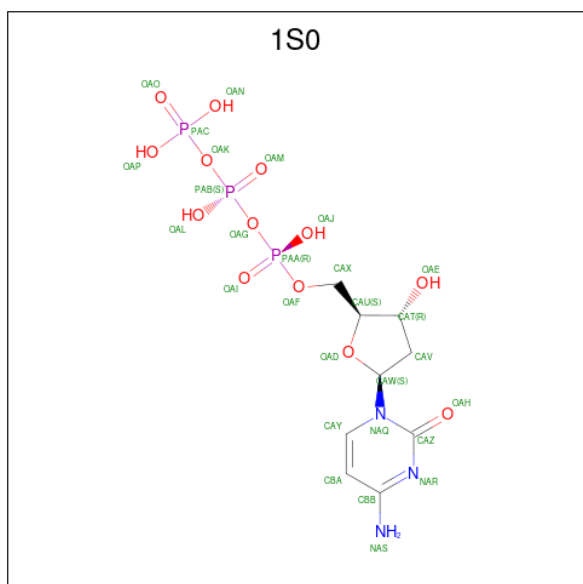
- Molecule 3 is a DNA chain called DNA (5'-D(*CP*AP*GP*TP*AP*C)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	6	Total	C	N	O	P	0	0	0
			119	58	23	33	5			
3	G	6	Total	C	N	O	P	0	0	0
			119	58	23	33	5			
3	K	6	Total	C	N	O	P	0	0	0
			119	58	23	33	5			
3	O	6	Total	C	N	O	P	0	0	0
			119	58	23	33	5			

- Molecule 4 is a DNA chain called DNA (5'-D(P*GP*CP*CP*G)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	4	Total 83	C 38	N 16	O 25	P 4	0	0	0
4	H	4	Total 83	C 38	N 16	O 25	P 4	0	0	0
4	L	4	Total 83	C 38	N 16	O 25	P 4	0	0	0
4	P	4	Total 83	C 38	N 16	O 25	P 4	0	0	0

- Molecule 5 is 4-amino-1-{2-deoxy-5-O-[(R)-hydroxy{[(S)-hydroxy(phosphonooxy)phosphoryl]oxy}phosphoryl]-beta-L-erythro-pentofuranosyl}pyrimidin-2(1H)-one (CCD ID: 1S0) (formula: C₉H₁₆N₃O₁₃P₃).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total 28	C 9	N 3	O 13	P 3	0	0
5	E	1	Total 28	C 9	N 3	O 13	P 3	0	0
5	I	1	Total 28	C 9	N 3	O 13	P 3	0	0
5	M	1	Total 28	C 9	N 3	O 13	P 3	0	0

- Molecule 6 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	5	Total	Ca	0	0
			5	5		

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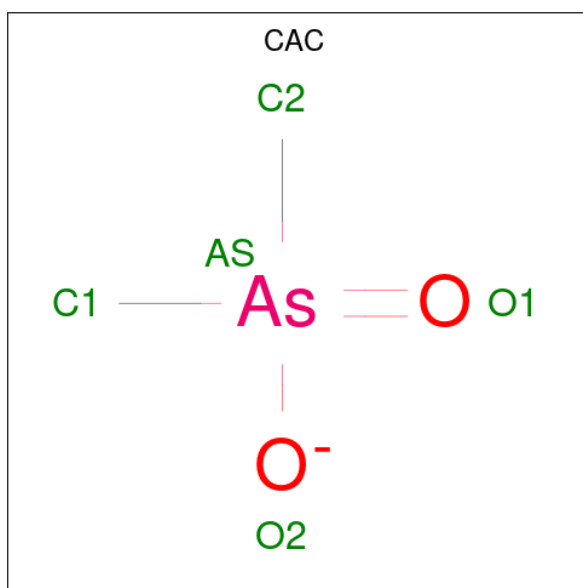
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	1	Total	Ca	0	0
			1	1		
6	E	5	Total	Ca	0	0
			5	5		
6	F	1	Total	Ca	0	0
			1	1		
6	I	3	Total	Ca	0	0
			3	3		
6	J	2	Total	Ca	0	0
			2	2		
6	L	1	Total	Ca	0	0
			1	1		
6	M	4	Total	Ca	0	0
			4	4		

- Molecule 7 is ACETATE ION (CCD ID: ACT) (formula: $C_2H_3O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	B	1	Total	C	O	0	0
			4	2	2		
7	F	1	Total	C	O	0	0
			4	2	2		
7	I	1	Total	C	O	0	0
			4	2	2		
7	M	1	Total	C	O	0	0
			4	2	2		

- Molecule 8 is CACODYLATE ION (CCD ID: CAC) (formula: $C_2H_6AsO_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
8	M	1	Total	As	C	O	0	0
			5	1	2	2		

- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	112	Total	O	0	0
			112	112		
9	B	21	Total	O	0	0
			21	21		
9	C	11	Total	O	0	0
			11	11		
9	E	64	Total	O	0	0
			64	64		
9	F	7	Total	O	0	0
			7	7		
9	G	4	Total	O	0	0
			4	4		
9	H	1	Total	O	0	0
			1	1		
9	I	66	Total	O	0	0
			66	66		
9	J	5	Total	O	0	0
			5	5		
9	K	14	Total	O	0	0
			14	14		

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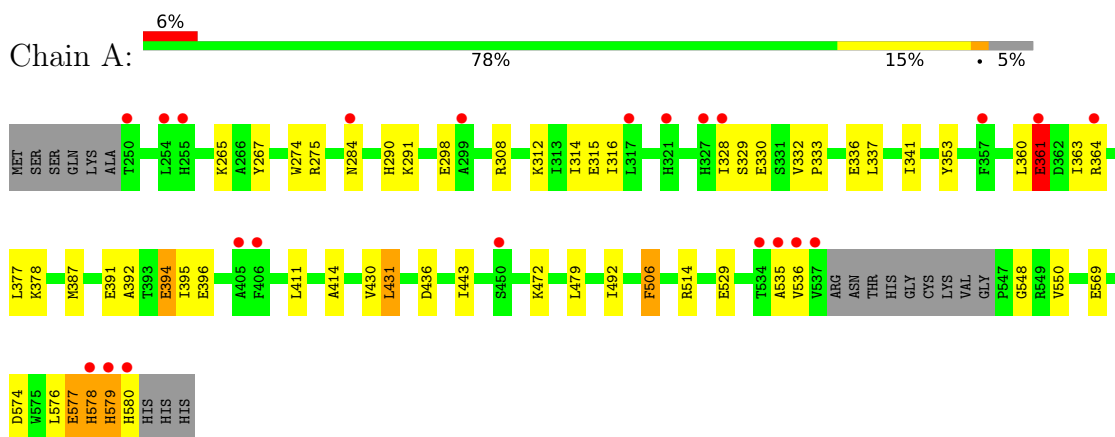
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	L	1	Total 1	O 1	0	0
9	M	105	Total 105	O 105	0	0
9	N	7	Total 7	O 7	0	0
9	O	2	Total 2	O 2	0	0
9	P	2	Total 2	O 2	0	0

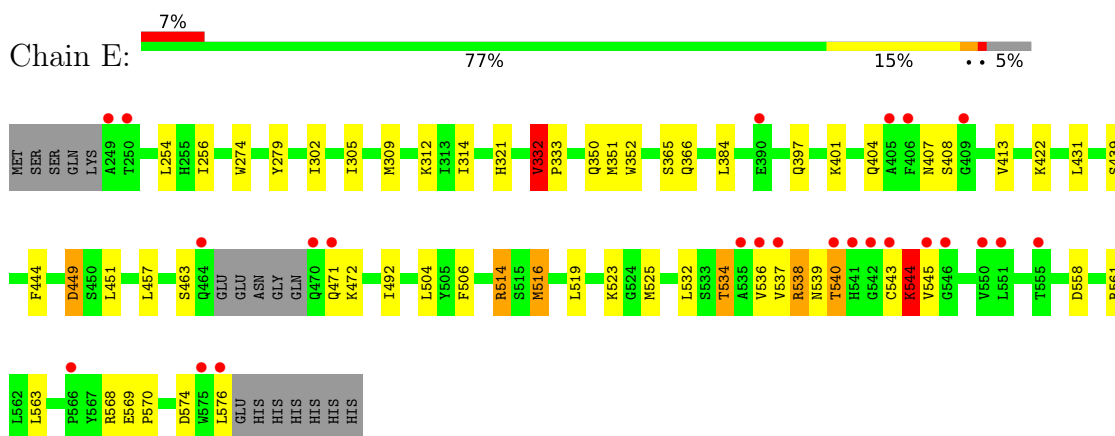
3 Residue-property plots [i](#)

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

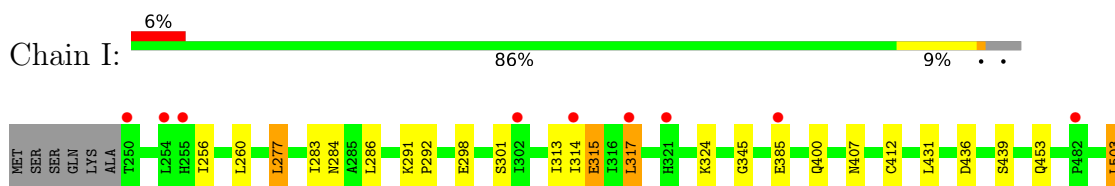
- Molecule 1: DNA polymerase lambda

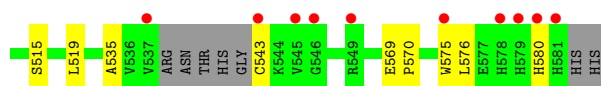


- Molecule 1: DNA polymerase lambda

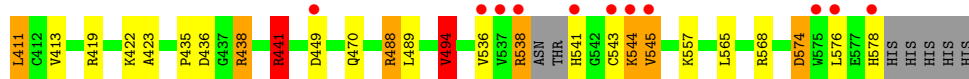
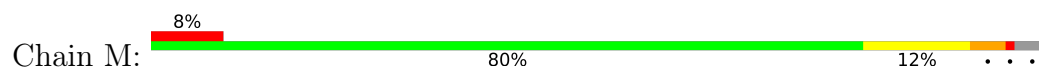


- Molecule 1: DNA polymerase lambda





- Molecule 1: DNA polymerase lambda



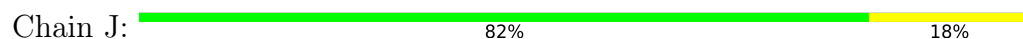
- Molecule 2: DNA (5'-D(*CP*GP*GP*CP*GP*GP*TP*AP*CP*TP*G)-3')



- Molecule 2: DNA (5'-D(*CP*GP*GP*CP*GP*GP*TP*AP*CP*TP*G)-3')



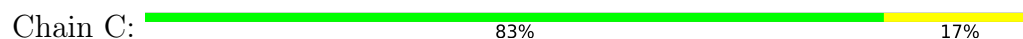
- Molecule 2: DNA (5'-D(*CP*GP*GP*CP*GP*GP*TP*AP*CP*TP*G)-3')



- Molecule 2: DNA (5'-D(*CP*GP*GP*CP*GP*GP*TP*AP*CP*TP*G)-3')



- Molecule 3: DNA (5'-D(*CP*AP*GP*TP*AP*C)-3')



- Molecule 3: DNA (5'-D(*CP*AP*GP*TP*AP*C)-3')

Chain G:  83% 17%



- Molecule 3: DNA (5'-D(*CP*AP*GP*TP*AP*C)-3')

Chain K:  83% 17%



- Molecule 3: DNA (5'-D(*CP*AP*GP*TP*AP*C)-3')

Chain O:  100%

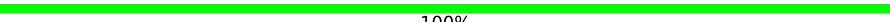
There are no outlier residues recorded for this chain.

- Molecule 4: DNA (5'-D(P*GP*CP*CP*G)-3')

Chain D:  50% 50%

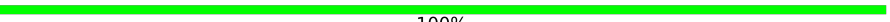


- Molecule 4: DNA (5'-D(P*GP*CP*CP*G)-3')

Chain H:  100%

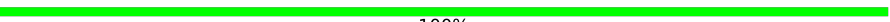
There are no outlier residues recorded for this chain.

- Molecule 4: DNA (5'-D(P*GP*CP*CP*G)-3')

Chain L:  100%

There are no outlier residues recorded for this chain.

- Molecule 4: DNA (5'-D(P*GP*CP*CP*G)-3')

Chain P:  100%

There are no outlier residues recorded for this chain.

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	194.57Å 97.81Å 105.07Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	36.60 – 2.15 36.60 – 2.15	Depositor EDS
% Data completeness (in resolution range)	98.1 (36.60-2.15) 96.7 (36.60-2.15)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.03 (at 2.16Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.200 , 0.250 (Not available) , 0.253	Depositor DCC
R_{free} test set	5379 reflections (4.91%)	wwPDB-VP
Wilson B-factor (Å ²)	33.5	Xtriage
Anisotropy	0.186	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 39.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	12538	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.88% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: ACT, 1S0, CA, CAC

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	1.22	3/2602 (0.1%)	1.12	9/3511 (0.3%)
1	E	1.04	2/2593 (0.1%)	1.08	2/3499 (0.1%)
1	I	1.06	1/2635 (0.0%)	1.01	1/3555 (0.0%)
1	M	1.19	3/2637 (0.1%)	1.12	6/3557 (0.2%)
2	B	0.50	0/252	0.94	0/388
2	F	0.40	0/252	1.07	1/388 (0.3%)
2	J	0.54	0/252	1.09	2/388 (0.5%)
2	N	0.54	0/252	0.92	1/388 (0.3%)
3	C	0.58	0/133	0.86	1/203 (0.5%)
3	G	0.55	0/133	1.12	2/203 (1.0%)
3	K	0.56	0/133	1.00	1/203 (0.5%)
3	O	0.50	0/133	0.99	0/203
4	D	0.57	0/92	0.96	0/138
4	H	0.47	0/92	0.87	0/138
4	L	0.42	0/92	0.81	0/138
4	P	0.45	0/92	0.93	0/138
All	All	1.06	9/12375 (0.1%)	1.07	26/17038 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	I	0	1
1	M	0	1
All	All	0	2

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	I	345	GLY	C-O	7.73	1.31	1.24
1	E	332	VAL	C-O	-6.53	1.19	1.24
1	A	332	VAL	CA-CB	6.13	1.57	1.54
1	M	289	PHE	CA-C	5.90	1.60	1.52
1	A	337	LEU	C-O	-5.85	1.17	1.24
1	A	341	ILE	C-O	-5.75	1.18	1.24
1	M	423	ALA	C-O	-5.38	1.17	1.24
1	M	441	ARG	CD-NE	-5.25	1.38	1.46
1	E	256	ILE	C-O	5.06	1.29	1.23

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	579	HIS	N-CA-C	-7.70	103.88	113.20
2	J	2	DG	C4'-C3'-O3'	-7.19	99.21	110.00
1	M	488	ARG	NE-CZ-NH2	-7.14	112.77	119.20
1	M	494	VAL	N-CA-CB	-7.06	101.32	111.21
2	F	8	DA	C2'-C3'-O3'	6.56	121.34	111.50
3	G	1	DC	C5'-C4'-O4'	6.37	118.96	109.40
2	N	1	DC	C5'-C4'-O4'	6.16	118.64	109.40
1	A	443	ILE	N-CA-C	6.06	116.80	110.62
1	M	419	ARG	N-CA-C	-6.00	105.44	112.89
1	E	274	TRP	N-CA-C	-5.87	105.22	112.90
1	E	449	ASP	N-CA-C	5.86	118.14	111.11
1	A	472	LYS	N-CA-C	5.84	118.06	109.24
2	J	1	DC	C5'-C4'-O4'	5.83	118.15	109.40
3	G	1	DC	C5'-C4'-C3'	5.49	123.14	114.90
3	K	3	DG	C4'-C3'-O3'	-5.45	101.82	110.00
1	A	284	ASN	N-CA-CB	5.30	117.69	110.01
1	I	277	LEU	CB-CA-C	-5.27	102.37	110.81
1	A	430	VAL	N-CA-C	5.26	115.53	108.17
1	A	316	ILE	CB-CA-C	-5.18	105.15	112.14
1	A	361	GLU	N-CA-C	-5.18	105.72	111.36
1	M	545	VAL	N-CA-C	5.17	120.10	109.34
3	C	3	DG	C2'-C3'-O3'	5.10	119.15	111.50
1	M	574	ASP	N-CA-C	-5.06	105.95	111.82
1	A	353	TYR	CA-C-N	5.05	127.99	120.31
1	A	353	TYR	C-N-CA	5.05	127.99	120.31
1	M	441	ARG	NE-CZ-NH2	-5.01	114.69	119.20

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	I	580	HIS	Peptide
1	M	290	HIS	Peptide

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	2548	0	2540	37	0
1	E	2541	0	2554	27	0
1	I	2580	0	2562	16	0
1	M	2584	0	2583	27	0
2	B	225	0	125	4	0
2	F	225	0	125	4	0
2	J	225	0	125	0	0
2	N	225	0	125	2	0
3	C	119	0	68	0	0
3	G	119	0	68	0	0
3	K	119	0	69	0	0
3	O	119	0	69	0	0
4	D	83	0	45	1	0
4	H	83	0	45	0	0
4	L	83	0	45	0	0
4	P	83	0	45	0	0
5	A	28	0	14	0	0
5	E	28	0	13	0	0
5	I	28	0	15	0	0
5	M	28	0	15	0	0
6	A	5	0	0	0	0
6	B	1	0	0	0	0
6	E	5	0	0	0	0
6	F	1	0	0	0	0
6	I	3	0	0	0	0
6	J	2	0	0	0	0
6	L	1	0	0	0	0
6	M	4	0	0	0	0
7	B	4	0	3	0	0
7	F	4	0	3	0	0
7	I	4	0	3	0	0
7	M	4	0	3	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
8	M	5	0	0	2	0
9	A	112	0	0	7	0
9	B	21	0	0	1	0
9	C	11	0	0	0	0
9	E	64	0	0	1	0
9	F	7	0	0	2	0
9	G	4	0	0	0	0
9	H	1	0	0	0	0
9	I	66	0	0	0	0
9	J	5	0	0	0	0
9	K	14	0	0	0	0
9	L	1	0	0	0	0
9	M	105	0	0	3	0
9	N	7	0	0	1	0
9	O	2	0	0	0	0
9	P	2	0	0	0	0
All	All	12538	0	11262	113	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 5.

All (113) 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:A:387:MET:HE1	1:A:395:ILE:HD12	1.20	1.16
1:E:543:CYS:HG	1:I:543:CYS:N	1.50	1.07
1:M:387:MET:HE1	1:M:395:ILE:HD12	1.37	1.05
1:A:387:MET:HE2	1:A:392:ALA:N	1.95	0.81
2:B:9:DC:OP1	9:B:208:HOH:O	2.00	0.77
1:A:387:MET:HE1	1:A:395:ILE:CD1	2.09	0.77
1:E:431:LEU:HD12	1:E:492:ILE:HG23	1.66	0.77
1:M:387:MET:HE2	1:M:392:ALA:N	2.00	0.76
2:N:7:DT:OP2	9:N:205:HOH:O	2.05	0.74
1:A:535:ALA:HB3	1:A:548:GLY:HA2	1.67	0.74
1:A:396:GLU:HG3	1:A:414:ALA:HB2	1.70	0.74
2:F:3:DG:N3	9:F:202:HOH:O	2.24	0.70
1:A:579:HIS:HB3	9:A:730:HOH:O	1.92	0.67
1:M:333:PRO:O	9:M:767:HOH:O	2.12	0.67
1:A:387:MET:HE3	1:A:391:GLU:HG2	1.76	0.67
1:M:422:LYS:NZ	1:M:574:ASP:OD2	2.29	0.66
1:M:438:ARG:O	1:M:441:ARG:HD3	1.97	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:351:MET:CE	1:E:352:TRP:CD1	2.81	0.64
1:E:543:CYS:SG	1:I:543:CYS:N	2.65	0.63
1:M:382:ASP:OD1	9:M:787:HOH:O	2.16	0.62
1:I:284:ASN:HB2	1:I:575:TRP:CE2	2.35	0.60
1:M:250:THR:O	1:M:251:ASN:CB	2.49	0.60
1:M:284:ASN:N	1:M:284:ASN:HD22	1.99	0.60
1:M:318:GLU:O	8:M:607:CAC:AS	2.80	0.59
1:A:333:PRO:O	9:A:725:HOH:O	2.17	0.59
1:A:387:MET:HE2	1:A:391:GLU:C	2.27	0.59
1:A:574:ASP:HB3	9:A:802:HOH:O	2.02	0.58
1:E:514:ARG:HD2	2:F:5:DG:OP2	2.04	0.57
1:M:291:LYS:HD3	1:M:292:PRO:O	2.05	0.57
2:B:1:DC:H2'	2:B:2:DG:C8	2.40	0.56
1:A:361:GLU:N	1:A:361:GLU:CD	2.63	0.56
1:A:394:GLU:HG3	1:A:479:LEU:HD11	1.88	0.56
1:A:431:LEU:C	1:A:431:LEU:HD23	2.31	0.56
1:M:438:ARG:O	1:M:441:ARG:CD	2.56	0.53
1:A:387:MET:HE3	1:A:391:GLU:CG	2.39	0.53
1:A:364:ARG:HG3	1:A:377:LEU:HD21	1.90	0.53
1:E:422:LYS:NZ	1:E:574:ASP:OD2	2.39	0.53
1:M:318:GLU:O	8:M:607:CAC:O2	2.27	0.53
1:A:308:ARG:O	1:A:312:LYS:HG3	2.10	0.52
1:I:436:ASP:C	1:I:436:ASP:OD1	2.53	0.52
1:E:407:ASN:ND2	1:E:439:SER:OG	2.44	0.51
1:M:290:HIS:CD2	1:M:291:LYS:HB2	2.47	0.50
1:A:387:MET:CE	1:A:391:GLU:HG2	2.42	0.50
1:M:255:HIS:H	1:M:255:HIS:CD2	2.30	0.50
1:I:314:ILE:HD12	1:I:315:GLU:N	2.27	0.50
1:M:557:LYS:HG3	9:M:776:HOH:O	2.12	0.50
1:E:558:ASP:OD1	1:E:561:ARG:NH2	2.45	0.49
1:M:357:PHE:O	1:M:358:ARG:NH1	2.37	0.49
1:M:436:ASP:C	1:M:436:ASP:OD1	2.55	0.49
1:E:538:ARG:HA	1:E:544:LYS:O	2.12	0.49
1:E:504:LEU:HD21	1:E:516:MET:CE	2.42	0.49
1:M:470:GLN:HG2	1:M:494:VAL:HG22	1.94	0.49
1:M:256:ILE:HD13	1:M:313:ILE:HG23	1.94	0.49
1:E:302:ILE:HB	1:E:305:ILE:HD12	1.95	0.49
1:A:330:GLU:O	1:A:333:PRO:HD2	2.13	0.49
1:A:291:LYS:NZ	1:A:298:GLU:OE2	2.45	0.48
1:I:569:GLU:O	1:I:570:PRO:C	2.57	0.48
1:E:279:TYR:OH	1:E:312:LYS:HE3	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:400:GLN:HG3	1:I:412:CYS:HB2	1.95	0.48
1:I:407:ASN:ND2	1:I:439:SER:OG	2.47	0.47
1:A:387:MET:HE2	1:A:392:ALA:CA	2.44	0.47
1:A:577:GLU:H	1:A:577:GLU:CD	2.23	0.46
1:A:265:LYS:HE3	1:A:336:GLU:HG3	1.96	0.46
1:A:314:ILE:O	1:A:315:GLU:C	2.58	0.46
2:N:9:DC:H2'	2:N:10:DT:H71	1.97	0.46
1:E:519:LEU:C	1:E:519:LEU:HD23	2.41	0.46
1:E:431:LEU:CD1	1:E:492:ILE:HG23	2.44	0.45
1:I:256:ILE:HD11	1:I:317:LEU:CD1	2.47	0.45
1:E:525:MET:CE	1:E:563:LEU:HD23	2.47	0.45
1:E:332:VAL:N	1:E:333:PRO:CD	2.80	0.45
1:A:579:HIS:CB	9:A:730:HOH:O	2.58	0.44
4:D:3:DC:H2''	4:D:4:DG:C8	2.52	0.44
1:I:515:SER:HB3	1:I:576:LEU:HD23	1.99	0.44
1:E:321:HIS:HB3	9:E:708:HOH:O	2.16	0.44
1:E:534:THR:O	1:E:536:VAL:HG23	2.18	0.44
1:I:291:LYS:HB2	1:I:292:PRO:HD2	1.99	0.44
1:M:538:ARG:HB3	1:M:538:ARG:CZ	2.47	0.44
1:A:267:TYR:CZ	1:A:275:ARG:HD3	2.52	0.44
1:A:360:LEU:O	1:A:363:ILE:HB	2.18	0.44
1:I:314:ILE:HD12	1:I:314:ILE:C	2.43	0.44
1:A:514:ARG:HD2	2:B:5:DG:OP2	2.18	0.43
2:F:1:DC:H2''	2:F:2:DG:H5'	1.99	0.43
1:A:514:ARG:HD3	2:B:5:DG:N7	2.34	0.43
1:E:523:LYS:NZ	1:E:563:LEU:O	2.48	0.43
1:E:534:THR:C	1:E:536:VAL:HG23	2.43	0.43
1:E:365:SER:OG	1:E:366:GLN:HG3	2.17	0.43
1:M:300:CYS:C	1:M:302:ILE:H	2.26	0.43
1:E:523:LYS:O	1:E:525:MET:HG3	2.18	0.42
1:M:541:HIS:O	1:M:541:HIS:ND1	2.52	0.42
1:I:431:LEU:HD23	1:I:503:LEU:HD13	2.01	0.42
1:A:274:TRP:HA	1:A:274:TRP:CE3	2.54	0.42
1:E:569:GLU:HG3	1:E:570:PRO:HD2	2.01	0.42
1:M:541:HIS:HA	1:M:543:CYS:SG	2.59	0.42
1:E:451:LEU:HB3	1:E:457:LEU:HG	2.02	0.42
1:E:525:MET:HE3	1:E:563:LEU:HD23	2.01	0.42
1:E:539:ASN:O	1:E:540:THR:C	2.62	0.42
1:I:286:LEU:HD11	1:I:313:ILE:HD11	2.00	0.41
1:A:578:HIS:C	1:A:580:HIS:N	2.78	0.41
2:F:9:DC:H5''	9:F:206:HOH:O	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:436:ASP:OD1	1:A:436:ASP:C	2.64	0.41
1:M:411:LEU:HD22	1:M:435:PRO:HD3	2.01	0.41
1:I:298:GLU:O	1:I:301:SER:OG	2.39	0.41
1:A:514:ARG:NH1	9:A:792:HOH:O	2.54	0.41
1:E:305:ILE:HG23	1:E:309:MET:HB3	2.01	0.41
1:A:492:ILE:HG23	1:A:506:PHE:CZ	2.56	0.41
1:I:283:ILE:O	1:I:284:ASN:C	2.64	0.40
1:M:288:SER:O	1:M:289:PHE:C	2.62	0.40
1:M:541:HIS:O	1:M:541:HIS:CG	2.73	0.40
1:A:328:ILE:O	1:A:329:SER:C	2.64	0.40
1:M:351:MET:CE	1:M:352:TRP:CD1	3.04	0.40
1:A:290:HIS:CE1	9:A:763:HOH:O	2.74	0.40
1:A:529:GLU:OE2	9:A:784:HOH:O	2.22	0.40
1:A:535:ALA:HB3	1:A:548:GLY:CA	2.45	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	318/340 (94%)	307 (96%)	9 (3%)	2 (1%)	21	15
1	E	319/340 (94%)	303 (95%)	12 (4%)	4 (1%)	9	5
1	I	323/340 (95%)	313 (97%)	9 (3%)	1 (0%)	36	34
1	M	324/340 (95%)	304 (94%)	14 (4%)	6 (2%)	6	2
All	All	1284/1360 (94%)	1227 (96%)	44 (3%)	13 (1%)	12	8

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	578	HIS

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Mol	Chain	Res	Type
1	M	251	ASN
1	I	535	ALA
1	M	301	SER
1	E	540	THR
1	E	544	LYS
1	E	545	VAL
1	M	329	SER
1	M	544	LYS
1	E	463	SER
1	M	290	HIS
1	A	536	VAL
1	M	545	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	272/287 (95%)	262 (96%)	10 (4%)	30	30
1	E	271/287 (94%)	247 (91%)	24 (9%)	9	5
1	I	274/287 (96%)	265 (97%)	9 (3%)	33	34
1	M	275/287 (96%)	252 (92%)	23 (8%)	10	6
All	All	1092/1148 (95%)	1026 (94%)	66 (6%)	17	13

All (66) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	361	GLU
1	A	378	LYS
1	A	394	GLU
1	A	411	LEU
1	A	431	LEU
1	A	506	PHE
1	A	550	VAL
1	A	569	GLU
1	A	576	LEU

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Mol	Chain	Res	Type
1	A	577	GLU
1	E	254	LEU
1	E	314	ILE
1	E	332	VAL
1	E	350	GLN
1	E	384	LEU
1	E	397	GLN
1	E	401	LYS
1	E	404	GLN
1	E	408	SER
1	E	413	VAL
1	E	444	PHE
1	E	449	ASP
1	E	471	GLN
1	E	472	LYS
1	E	506	PHE
1	E	514	ARG
1	E	516	MET
1	E	532	LEU
1	E	534	THR
1	E	537	VAL
1	E	538	ARG
1	E	544	LYS
1	E	568	ARG
1	E	576	LEU
1	I	260	LEU
1	I	277	LEU
1	I	315	GLU
1	I	317	LEU
1	I	324	LYS
1	I	385	GLU
1	I	453	GLN
1	I	503	LEU
1	I	519	LEU
1	M	250	THR
1	M	284	ASN
1	M	286	LEU
1	M	302	ILE
1	M	314	ILE
1	M	317	LEU
1	M	384	LEU
1	M	385	GLU

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Mol	Chain	Res	Type
1	M	411	LEU
1	M	413	VAL
1	M	438	ARG
1	M	441	ARG
1	M	449	ASP
1	M	488	ARG
1	M	489	LEU
1	M	494	VAL
1	M	536	VAL
1	M	538	ARG
1	M	544	LYS
1	M	565	LEU
1	M	568	ARG
1	M	576	LEU
1	M	578	HIS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (32) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	255	HIS
1	A	270	GLN
1	A	284	ASN
1	A	340	ASN
1	A	372	GLN
1	A	373	GLN
1	A	397	GLN
1	A	404	GLN
1	A	471	GLN
1	A	511	HIS
1	A	580	HIS
1	E	270	GLN
1	E	290	HIS
1	E	327	HIS
1	E	372	GLN
1	E	404	GLN
1	E	541	HIS
1	I	284	ASN
1	I	355	GLN
1	I	372	GLN
1	I	397	GLN
1	I	404	GLN
1	I	453	GLN

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Mol	Chain	Res	Type
1	I	471	GLN
1	M	252	HIS
1	M	255	HIS
1	M	284	ASN
1	M	290	HIS
1	M	350	GLN
1	M	372	GLN
1	M	379	HIS
1	M	486	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 31 ligands modelled in this entry, 22 are monoatomic - leaving 9 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	$\# Z > 2$
5	1S0	M	602	6	28,29,29	1.30	2 (7%)	41,45,45	1.79	8 (19%)
5	1S0	I	601	6	28,29,29	2.26	5 (17%)	41,45,45	1.53	4 (9%)
5	1S0	E	601	6	28,29,29	1.99	5 (17%)	41,45,45	1.75	5 (12%)
5	1S0	A	601	6	28,29,29	1.79	5 (17%)	41,45,45	1.37	5 (12%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
8	CAC	M	607	-	2,4,4	3.08	2 (100%)	4,6,6	2.87	2 (50%)
7	ACT	F	102	-	3,3,3	0.75	0	3,3,3	0.98	0
7	ACT	I	605	-	3,3,3	0.72	0	3,3,3	1.31	0
7	ACT	B	102	-	3,3,3	0.85	0	3,3,3	0.73	0
7	ACT	M	601	-	3,3,3	0.88	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
5	1S0	M	602	6	-	2/22/34/34	0/2/2/2
5	1S0	E	601	6	-	5/22/34/34	0/2/2/2
5	1S0	A	601	6	-	4/22/34/34	0/2/2/2
5	1S0	I	601	6	-	1/22/34/34	0/2/2/2

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	I	601	1S0	OAH-CAZ	7.80	1.38	1.23
5	E	601	1S0	PAA-OAG	7.42	1.67	1.59
5	I	601	1S0	PAA-OAG	5.71	1.65	1.59
5	A	601	1S0	OAH-CAZ	5.42	1.33	1.23
5	A	601	1S0	PAA-OAG	5.32	1.65	1.59
5	M	602	1S0	OAH-CAZ	4.96	1.33	1.23
5	I	601	1S0	PAB-OAG	4.72	1.64	1.59
5	E	601	1S0	OAH-CAZ	4.32	1.31	1.23
5	I	601	1S0	PAB-OAK	3.74	1.63	1.59
5	E	601	1S0	PAB-OAG	3.16	1.62	1.59
8	M	607	CAC	AS-C2	3.10	1.97	1.90
8	M	607	CAC	AS-C1	3.07	1.97	1.90
5	M	602	1S0	CAY-CBA	2.72	1.41	1.35
5	E	601	1S0	PAB-OAK	2.46	1.62	1.59
5	A	601	1S0	CAZ-NAQ	-2.44	1.35	1.40
5	I	601	1S0	CAY-CBA	2.33	1.40	1.35
5	E	601	1S0	CAZ-NAQ	-2.27	1.35	1.40
5	A	601	1S0	CAZ-NAR	-2.12	1.32	1.36
5	A	601	1S0	OAE-CAT	-2.07	1.39	1.43

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	E	601	1S0	OAJ-PAA-OAG	6.15	123.90	107.27
5	M	602	1S0	OAH-CAZ-NAR	-5.63	113.46	122.33
5	I	601	1S0	OAH-CAZ-NAR	-5.48	113.69	122.33
5	M	602	1S0	OAD-CAW-NAQ	4.67	116.14	107.86
5	E	601	1S0	OAH-CAZ-NAR	-4.66	114.98	122.33
5	A	601	1S0	OAH-CAZ-NAR	-4.56	115.15	122.33
8	M	607	CAC	O1-AS-C2	-4.53	105.87	111.50
5	M	602	1S0	NAQ-CAZ-NAR	3.73	125.27	118.80
5	I	601	1S0	OAH-CAZ-NAQ	3.54	125.84	118.90
8	M	607	CAC	O2-AS-C2	3.34	113.93	105.84
5	A	601	1S0	OAG-PAA-OAI	-3.05	101.52	110.70
5	A	601	1S0	NAQ-CAZ-NAR	2.98	123.97	118.80
5	I	601	1S0	CAW-NAQ-CAZ	2.95	122.88	117.83
5	E	601	1S0	NAQ-CAZ-NAR	2.68	123.46	118.80
5	E	601	1S0	OAL-PAB-OAM	2.55	124.33	112.44
5	A	601	1S0	OAJ-PAA-OAI	2.53	124.19	112.44
5	E	601	1S0	CAW-NAQ-CAZ	2.52	122.15	117.83
5	M	602	1S0	CAY-CBA-CBB	2.50	121.63	117.53
5	M	602	1S0	CBA-CAY-NAQ	-2.41	117.92	121.84
5	M	602	1S0	OAL-PAB-OAM	2.29	123.11	112.44
5	A	601	1S0	OAP-PAC-OAO	2.22	119.50	110.83
5	M	602	1S0	OAP-PAC-OAK	-2.18	97.31	104.64
5	M	602	1S0	OAL-PAB-OAK	2.16	113.10	107.27
5	I	601	1S0	OAN-PAC-OAP	2.14	115.84	107.80

There are no chirality outliers.

All (12) torsion outliers are listed below:

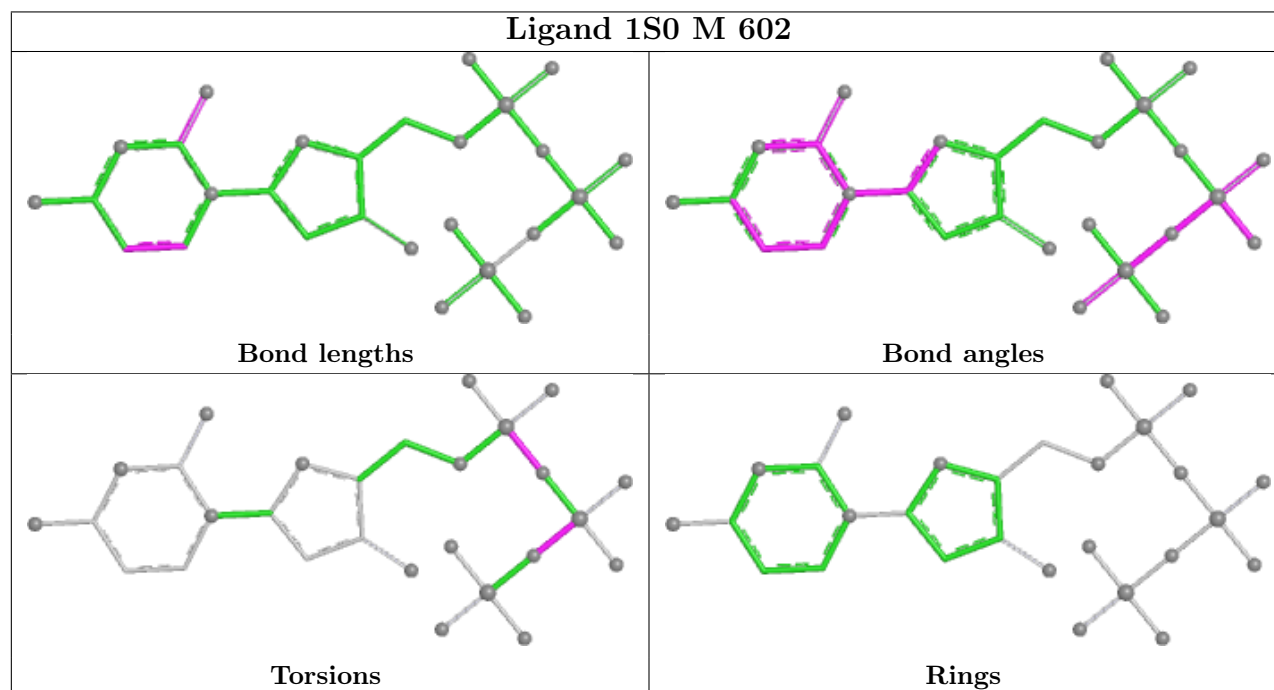
Mol	Chain	Res	Type	Atoms
5	A	601	1S0	PAB-OAK-PAC-OAP
5	E	601	1S0	PAB-OAK-PAC-OAP
5	I	601	1S0	PAB-OAG-PAA-OAF
5	M	602	1S0	PAB-OAG-PAA-OAF
5	E	601	1S0	PAA-OAG-PAB-OAM
5	A	601	1S0	PAB-OAK-PAC-OAO
5	A	601	1S0	PAA-OAG-PAB-OAM
5	E	601	1S0	PAA-OAG-PAB-OAL
5	E	601	1S0	PAB-OAK-PAC-OAO
5	A	601	1S0	PAA-OAG-PAB-OAL
5	E	601	1S0	PAB-OAG-PAA-OAJ
5	M	602	1S0	PAC-OAK-PAB-OAL

There are no ring outliers.

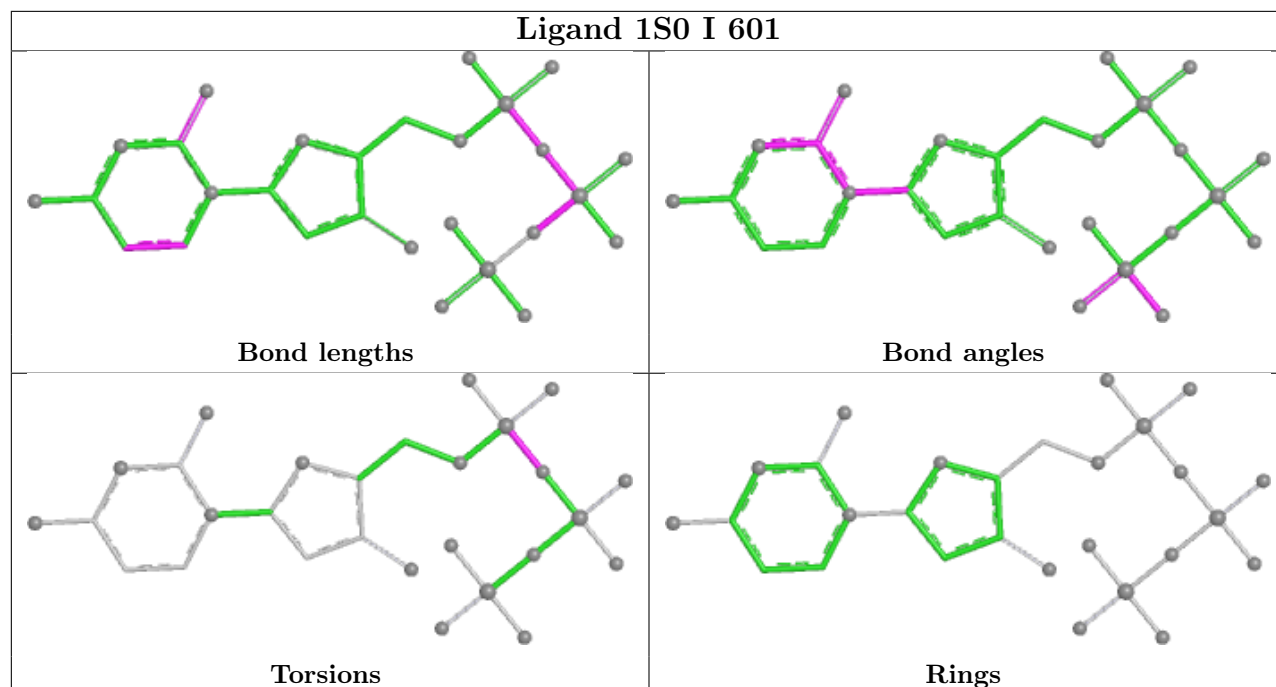
1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
8	M	607	CAC	2	0

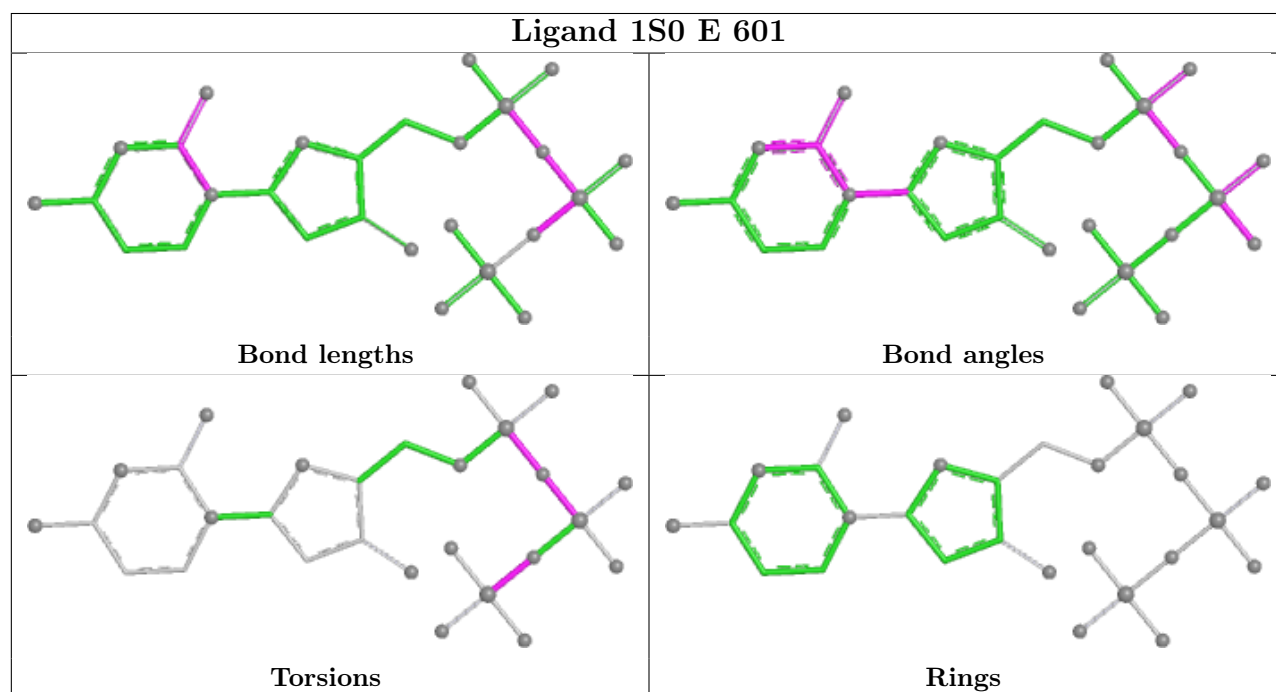
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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.

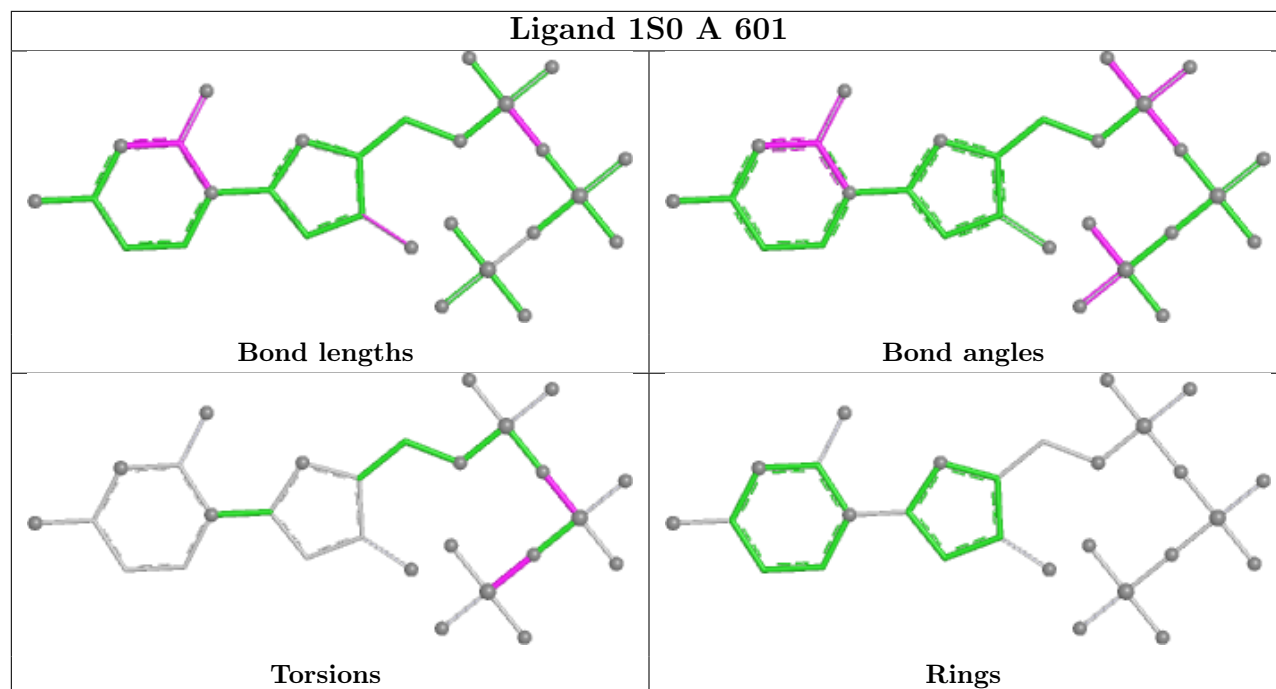


Ligand 1S0 I 601



Ligand 1S0 E 601





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	322/340 (94%)	0.31	22 (6%) 23 26	17, 34, 60, 112	0
1	E	323/340 (95%)	0.57	24 (7%) 20 23	20, 40, 70, 118	0
1	I	327/340 (96%)	0.52	19 (5%) 29 32	22, 42, 70, 104	0
1	M	328/340 (96%)	0.45	26 (7%) 18 20	17, 38, 60, 104	0
2	B	11/11 (100%)	-0.67	0 100 100	21, 26, 44, 44	0
2	F	11/11 (100%)	0.34	0 100 100	38, 43, 54, 62	0
2	J	11/11 (100%)	-0.17	0 100 100	36, 40, 43, 50	0
2	N	11/11 (100%)	-0.38	0 100 100	24, 32, 42, 50	0
3	C	6/6 (100%)	-0.77	0 100 100	19, 23, 38, 43	0
3	G	6/6 (100%)	0.00	0 100 100	32, 35, 62, 73	0
3	K	6/6 (100%)	-0.32	0 100 100	25, 29, 55, 57	0
3	O	6/6 (100%)	-0.29	0 100 100	28, 32, 52, 59	0
4	D	4/4 (100%)	-0.38	0 100 100	33, 35, 38, 39	0
4	H	4/4 (100%)	-0.40	0 100 100	32, 34, 38, 44	0
4	L	4/4 (100%)	0.39	0 100 100	44, 46, 51, 51	0
4	P	4/4 (100%)	-0.16	0 100 100	34, 38, 45, 47	0
All	All	1384/1444 (95%)	0.42	91 (6%) 24 27	17, 38, 67, 118	0

All (91) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	M	575	TRP	5.3
1	A	537	VAL	4.8
1	I	578	HIS	4.5
1	M	545	VAL	4.5
1	A	255	HIS	4.0

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Mol	Chain	Res	Type	RSRZ
1	M	537	VAL	3.9
1	E	540	THR	3.9
1	A	580	HIS	3.8
1	M	576	LEU	3.7
1	I	302	ILE	3.7
1	E	541	HIS	3.6
1	I	581	HIS	3.6
1	E	537	VAL	3.6
1	M	302	ILE	3.6
1	A	535	ALA	3.5
1	M	449	ASP	3.5
1	I	579	HIS	3.5
1	E	545	VAL	3.4
1	M	544	LYS	3.4
1	E	249	ALA	3.4
1	M	578	HIS	3.3
1	E	576	LEU	3.3
1	M	289	PHE	3.2
1	I	546	GLY	3.2
1	I	575	TRP	3.2
1	E	536	VAL	3.1
1	E	575	TRP	3.1
1	A	536	VAL	3.1
1	M	536	VAL	3.1
1	M	541	HIS	3.1
1	I	537	VAL	3.0
1	A	284	ASN	3.0
1	M	250	THR	3.0
1	M	303	PRO	2.9
1	I	250	THR	2.9
1	A	579	HIS	2.9
1	A	250	THR	2.9
1	A	534	THR	2.9
1	A	327	HIS	2.8
1	M	385	GLU	2.8
1	E	550	VAL	2.8
1	A	406	PHE	2.7
1	M	538	ARG	2.7
1	I	543	CYS	2.7
1	E	470	GLN	2.7
1	E	555	THR	2.7
1	E	409	GLY	2.7

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Mol	Chain	Res	Type	RSRZ
1	E	542	GLY	2.7
1	I	385	GLU	2.6
1	M	288	SER	2.6
1	I	255	HIS	2.5
1	A	578	HIS	2.5
1	E	250	THR	2.5
1	M	249	ALA	2.5
1	A	357	PHE	2.5
1	A	328	ILE	2.5
1	E	471	GLN	2.5
1	M	384	LEU	2.4
1	M	543	CYS	2.4
1	E	405	ALA	2.4
1	M	285	ALA	2.4
1	I	317	LEU	2.4
1	I	254	LEU	2.4
1	I	482	PRO	2.4
1	E	535	ALA	2.3
1	A	317	LEU	2.3
1	E	551	LEU	2.3
1	I	321	HIS	2.3
1	I	580	HIS	2.3
1	M	362	ASP	2.3
1	E	406	PHE	2.3
1	I	545	VAL	2.3
1	A	364	ARG	2.2
1	A	299	ALA	2.2
1	M	351	MET	2.2
1	I	314	ILE	2.2
1	E	390	GLU	2.2
1	A	321	HIS	2.2
1	M	286	LEU	2.2
1	I	549	ARG	2.2
1	E	543	CYS	2.1
1	A	361	GLU	2.1
1	A	254	LEU	2.1
1	A	405	ALA	2.1
1	E	464	GLN	2.0
1	M	283	ILE	2.0
1	A	450	SER	2.0
1	M	280	ALA	2.0
1	M	282	ALA	2.0

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Mol	Chain	Res	Type	RSRZ
1	E	566	PRO	2.0
1	E	546	GLY	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. 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.

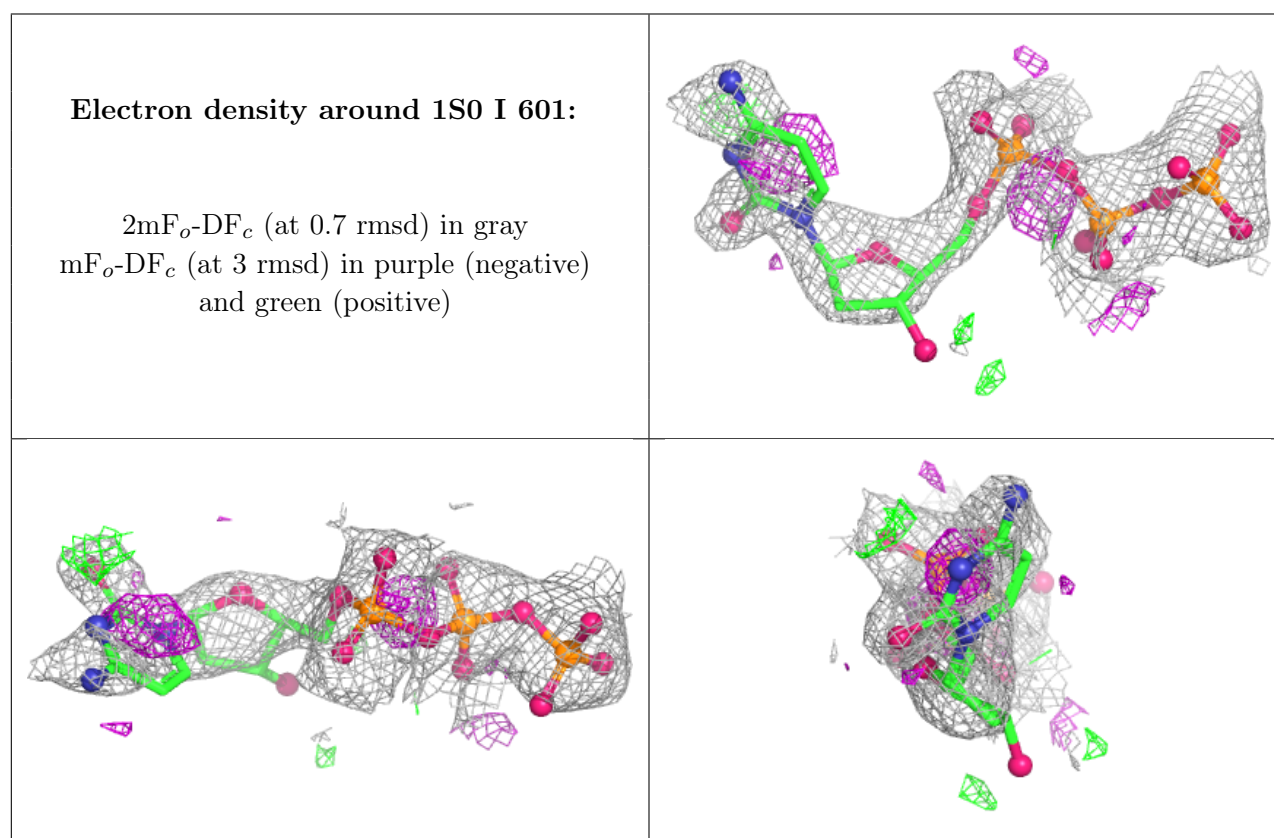
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
7	ACT	B	102	4/4	0.64	0.28	60,62,63,71	0
6	CA	E	604	1/1	0.74	0.30	78,78,78,78	0
7	ACT	F	102	4/4	0.79	0.14	45,51,53,57	0
6	CA	I	602	1/1	0.81	0.17	69,69,69,69	0
6	CA	M	606	1/1	0.82	0.25	72,72,72,72	0
7	ACT	M	601	4/4	0.83	0.27	54,58,61,64	0
6	CA	J	102	1/1	0.84	0.14	81,81,81,81	0
8	CAC	M	607	5/5	0.84	0.19	96,106,107,123	0
7	ACT	I	605	4/4	0.85	0.17	46,53,55,55	0
6	CA	M	604	1/1	0.85	0.17	72,72,72,72	0
5	ISO	I	601	28/28	0.85	0.17	51,76,84,84	0
6	CA	E	605	1/1	0.88	0.13	56,56,56,56	0
6	CA	J	101	1/1	0.90	0.11	67,67,67,67	0
6	CA	A	605	1/1	0.91	0.24	57,57,57,57	0
5	ISO	M	602	28/28	0.91	0.10	31,47,58,61	0
6	CA	I	603	1/1	0.91	0.15	77,77,77,77	0
6	CA	M	603	1/1	0.93	0.12	67,67,67,67	0
6	CA	L	101	1/1	0.95	0.06	78,78,78,78	0
6	CA	A	604	1/1	0.95	0.25	34,34,34,34	0
6	CA	F	101	1/1	0.95	0.06	56,56,56,56	0
6	CA	B	101	1/1	0.96	0.06	59,59,59,59	0

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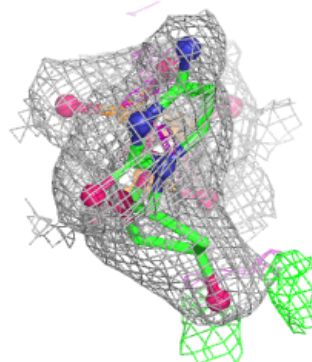
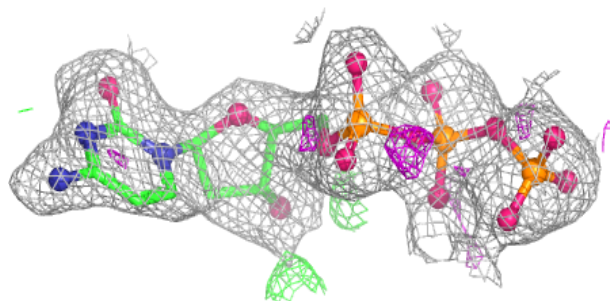
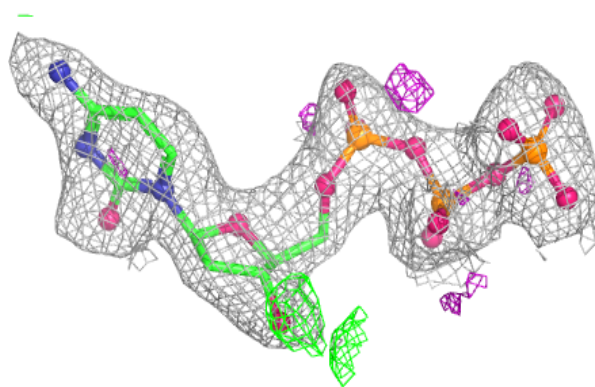
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
6	CA	E	606	1/1	0.96	0.31	41,41,41,41	0
6	CA	M	605	1/1	0.96	0.21	40,40,40,40	0
6	CA	A	603	1/1	0.97	0.05	20,20,20,20	0
5	1S0	E	601	28/28	0.97	0.06	23,30,39,43	0
6	CA	E	603	1/1	0.97	0.10	37,37,37,37	0
6	CA	I	604	1/1	0.97	0.03	20,20,20,20	1
6	CA	A	606	1/1	0.98	0.11	42,42,42,42	0
6	CA	E	602	1/1	0.99	0.05	34,34,34,34	0
6	CA	A	602	1/1	0.99	0.03	20,20,20,20	0
5	1S0	A	601	28/28	0.99	0.04	14,18,21,23	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.

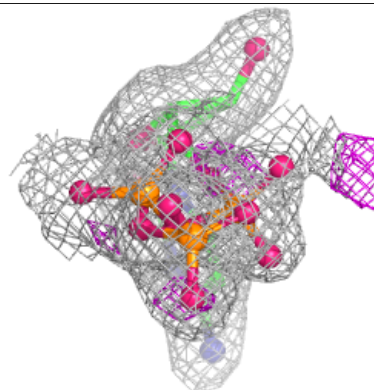
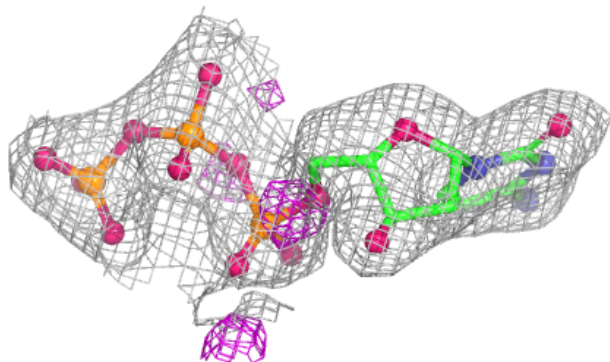
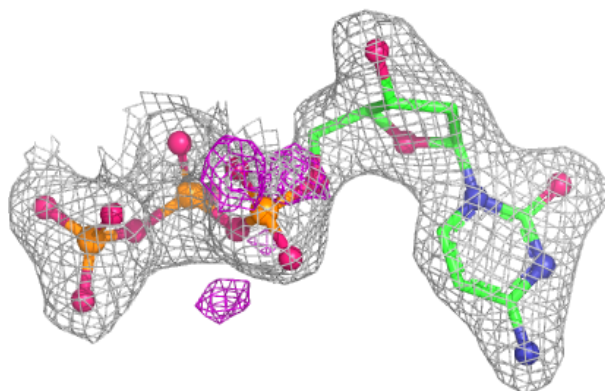


Electron density around 1S0 M 602:

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

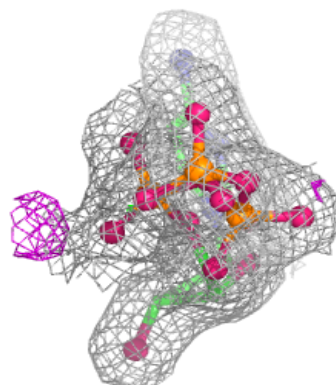
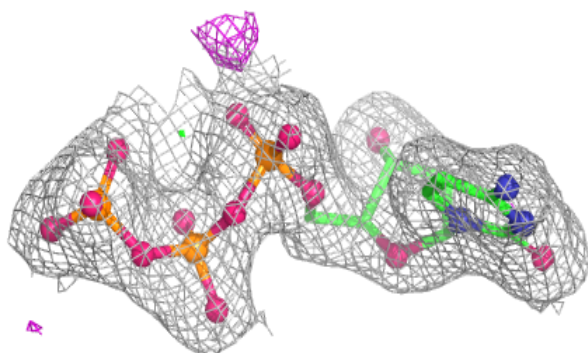
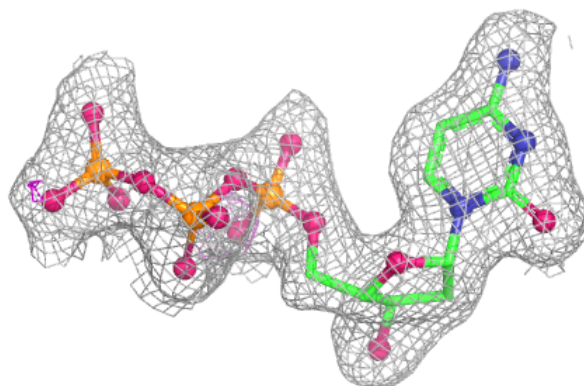
**Electron density around 1S0 E 601:**

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



Electron density around 1S0 A 601:

$2mF_o - DF_c$ (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.