



Full wwPDB X-ray Structure Validation Report ⓘ

May 22, 2020 – 01:29 am BST

PDB ID : 4QRH
Title : Molecular mechanism and evolution of guanylate kinase regulation by (p)ppGpp
Authors : Liu, K.; Krasny, K.; Myers, A.R.; Satyshur, K.A.; Keck, J.L.; Wnag, J.D.
Deposited on : 2014-07-01
Resolution : 1.65 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

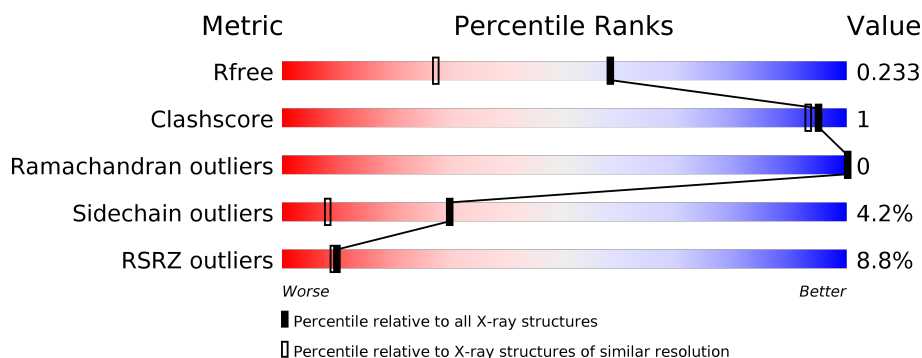
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.65 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}	130704	1827 (1.66-1.66)
Clashscore	141614	1931 (1.66-1.66)
Ramachandran outliers	138981	1891 (1.66-1.66)
Sidechain outliers	138945	1891 (1.66-1.66)
RSRZ outliers	127900	1791 (1.66-1.66)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria 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	210	<div> <div>7%</div> <div>81%</div> <div>14%</div> </div>
1	B	210	<div> <div>8%</div> <div>83%</div> <div>5%</div> <div>11%</div> </div>
1	C	210	<div> <div>9%</div> <div>82%</div> <div>6%</div> <div>12%</div> </div>
1	D	210	<div> <div>8%</div> <div>81%</div> <div>6%</div> <div>12%</div> </div>

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 6870 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 Guanylate kinase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	180	Total	C	N	O	S	0	4	0
			1493	952	249	283	9			
1	B	186	Total	C	N	O	S	0	7	0
			1564	995	263	298	8			
1	C	185	Total	C	N	O	S	0	3	0
			1523	972	251	292	8			
1	D	184	Total	C	N	O	S	0	2	0
			1513	962	254	290	7			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	SER	-	EXPRESSION TAG	UNP W6E076
A	-1	ASN	-	EXPRESSION TAG	UNP W6E076
A	0	ALA	-	EXPRESSION TAG	UNP W6E076
B	-2	SER	-	EXPRESSION TAG	UNP W6E076
B	-1	ASN	-	EXPRESSION TAG	UNP W6E076
B	0	ALA	-	EXPRESSION TAG	UNP W6E076
C	-2	SER	-	EXPRESSION TAG	UNP W6E076
C	-1	ASN	-	EXPRESSION TAG	UNP W6E076
C	0	ALA	-	EXPRESSION TAG	UNP W6E076
D	-2	SER	-	EXPRESSION TAG	UNP W6E076
D	-1	ASN	-	EXPRESSION TAG	UNP W6E076
D	0	ALA	-	EXPRESSION TAG	UNP W6E076

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		

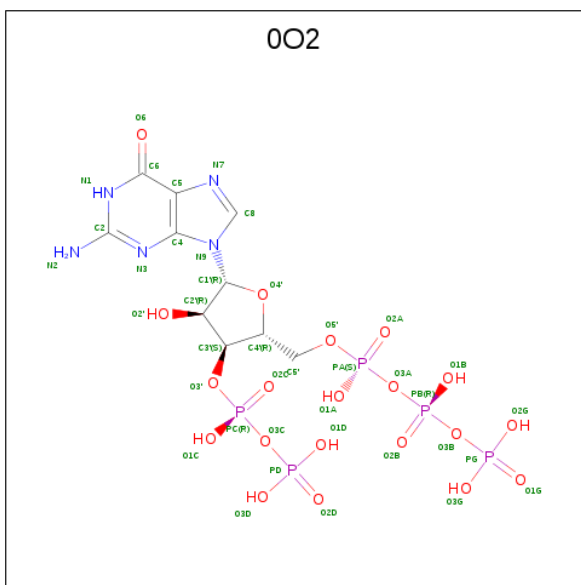
- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Mg	0	0
			1	1		
3	D	1	Total	Mg	0	0
			1	1		
3	C	1	Total	Mg	0	0
			1	1		

- Molecule 4 is POTASSIUM ION (three-letter code: K) (formula: K).

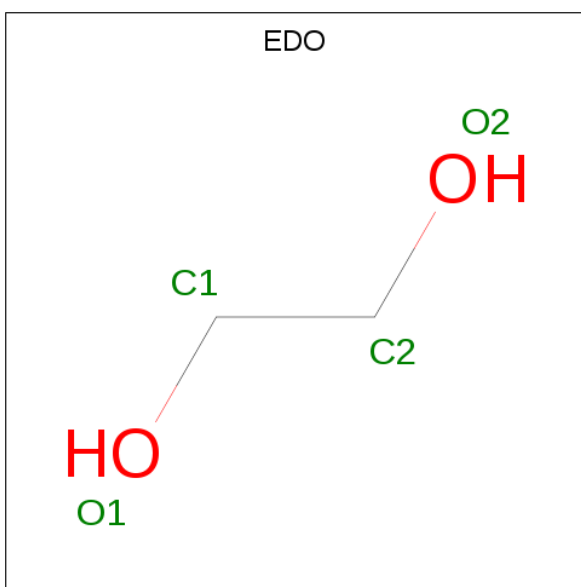
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	K	0	0
			1	1		
4	D	1	Total	K	0	0
			1	1		
4	C	1	Total	K	0	0
			1	1		

- Molecule 5 is guanosine 5'-(tetrahydrogen triphosphate) 3'-(trihydrogen diphosphate) (three-letter code: 0O2) (formula: C₁₀H₁₈N₅O₂₀P₅).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	B	1	Total 40	C 10	N 5	O 20	P 5	0	0
5	C	1	Total 40	C 10	N 5	O 20	P 5	0	0
5	D	1	Total 40	C 10	N 5	O 20	P 5	0	0

- Molecule 6 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: $\text{C}_2\text{H}_6\text{O}_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	D	1	Total	C	O	0	0
			4	2	2		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	D	1	Total	C	O	0	0
			4	2	2		
6	D	1	Total	C	O	0	0
			4	2	2		

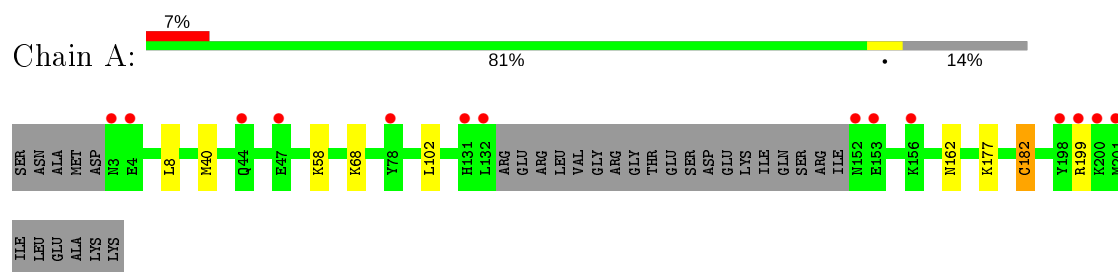
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	151	Total	O	0	0
			151	151		
7	B	184	Total	O	0	0
			184	184		
7	C	161	Total	O	0	0
			161	161		
7	D	133	Total	O	0	0
			133	133		

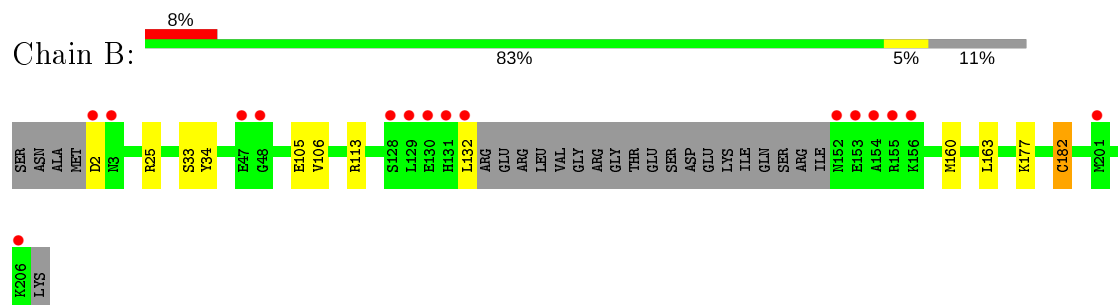
3 Residue-property plots [i](#)

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

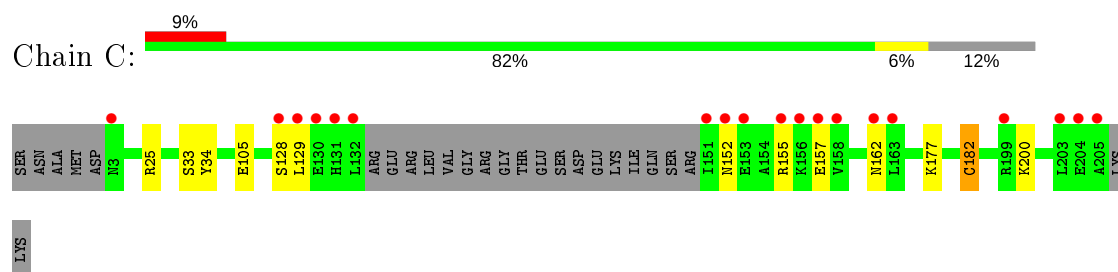
• Molecule 1: Guanylate kinase



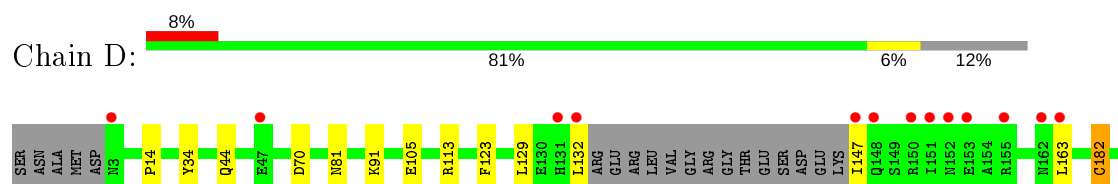
• Molecule 1: Guanylate kinase



• Molecule 1: Guanylate kinase



• Molecule 1: Guanylate kinase



HIS7	MET
Y198	ILE
H199	LEU
K200	GLU
	ALA
	LYS
	LYS

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	69.49 Å 93.96 Å 85.26 Å 90.00° 111.09° 90.00°	Depositor
Resolution (Å)	39.26 – 1.65 39.26 – 1.64	Depositor EDS
% Data completeness (in resolution range)	99.3 (39.26-1.65) 99.3 (39.26-1.64)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.62	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.09 (at 1.64 Å)	Xtriage
Refinement program	REFMAC 5.8.0071	Depositor
R, R_{free}	0.201 , 0.227 0.212 , 0.233	Depositor DCC
R_{free} test set	6171 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	19.8	Xtriage
Anisotropy	0.020	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 47.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.013 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6870	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.26% 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: SO4, MG, 0O2, K, EDO

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	0.35	0/1529	0.59	0/2054
1	B	0.35	0/1606	0.58	0/2156
1	C	0.35	0/1559	0.59	0/2096
1	D	0.35	0/1543	0.58	0/2074
All	All	0.35	0/6237	0.59	0/8380

There are no bond length outliers.

There are no bond angle outliers.

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	1493	0	1481	3	0
1	B	1564	0	1554	4	0
1	C	1523	0	1507	4	0
1	D	1513	0	1492	6	0
2	A	5	0	0	0	0
2	C	5	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	1	0	0	0	0
4	C	1	0	0	0	0
4	D	1	0	0	0	0
5	B	40	0	16	0	0
5	C	40	0	17	0	0
5	D	40	0	15	0	0
6	D	12	0	18	3	0
7	A	151	0	0	0	0
7	B	184	0	0	0	0
7	C	161	0	0	1	0
7	D	133	0	0	0	0
All	All	6870	0	6100	15	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 1.

All (15) 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:40[B]:MET:SD	1:A:58:LYS:HG3	2.37	0.65
1:D:44:GLN:HB2	6:D:305:EDO:HO2	1.72	0.54
1:C:129:LEU:HD21	1:C:157:GLU:HG2	1.90	0.53
1:C:25:ARG:HG3	1:C:177:LYS:HD3	1.91	0.52
1:A:182[B]:CYS:SG	1:B:182[B]:CYS:O	2.68	0.52
1:D:44:GLN:HB2	6:D:305:EDO:O2	2.10	0.52
1:D:113:ARG:NH1	1:D:163:LEU:O	2.40	0.47
1:D:81:ASN:ND2	6:D:305:EDO:H12	2.31	0.46
1:B:106:VAL:HG11	1:B:160:MET:HE3	1.98	0.46
1:C:182[B]:CYS:SG	1:D:182[B]:CYS:O	2.75	0.45
1:B:113[B]:ARG:NH2	1:B:163:LEU:O	2.47	0.44
1:B:25:ARG:HG3	1:B:177:LYS:HD3	2.00	0.44
1:D:14:PRO:HD3	1:D:123:PHE:CZ	2.55	0.41
1:A:8:LEU:HD11	1:A:102:LEU:HG	2.02	0.41
1:C:33:SER:HB2	7:C:560:HOH:O	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	180/210 (86%)	177 (98%)	3 (2%)	0	100	100
1	B	189/210 (90%)	187 (99%)	2 (1%)	0	100	100
1	C	184/210 (88%)	180 (98%)	4 (2%)	0	100	100
1	D	182/210 (87%)	179 (98%)	3 (2%)	0	100	100
All	All	735/840 (88%)	723 (98%)	12 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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	163/185 (88%)	157 (96%)	6 (4%)	34	10
1	B	171/185 (92%)	164 (96%)	7 (4%)	30	8
1	C	166/185 (90%)	157 (95%)	9 (5%)	22	4
1	D	165/185 (89%)	156 (94%)	9 (6%)	21	4
All	All	665/740 (90%)	634 (95%)	31 (5%)	30	6

All (31) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	68	LYS
1	A	162	ASN

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Mol	Chain	Res	Type
1	A	177	LYS
1	A	182[A]	CYS
1	A	182[B]	CYS
1	A	199	ARG
1	B	2	ASP
1	B	33	SER
1	B	34	TYR
1	B	105	GLU
1	B	132	LEU
1	B	182[A]	CYS
1	B	182[B]	CYS
1	C	34	TYR
1	C	105	GLU
1	C	128	SER
1	C	152	ASN
1	C	155	ARG
1	C	162	ASN
1	C	182[A]	CYS
1	C	182[B]	CYS
1	C	200	LYS
1	D	34	TYR
1	D	70	ASP
1	D	91	LYS
1	D	105	GLU
1	D	129	LEU
1	D	132	LEU
1	D	147	ILE
1	D	182[A]	CYS
1	D	182[B]	CYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	71	GLN
1	A	152	ASN
1	B	71	GLN
1	C	181	GLN
1	D	81	ASN
1	D	170	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 14 ligands modelled in this entry, 6 are monoatomic - leaving 8 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	EDO	D	305	-	3,3,3	0.51	0	2,2,2	0.55	0
6	EDO	D	304	-	3,3,3	0.48	0	2,2,2	0.30	0
5	0O2	C	304	3	32,42,42	1.65	5 (15%)	45,68,68	1.68	11 (24%)
2	SO4	A	301	-	4,4,4	0.29	0	6,6,6	0.32	0
5	0O2	D	306	3	32,42,42	1.63	5 (15%)	45,68,68	1.57	10 (22%)
6	EDO	D	303	-	3,3,3	0.58	0	2,2,2	0.11	0
2	SO4	C	303	-	4,4,4	0.26	0	6,6,6	0.36	0
5	0O2	B	303	3	32,42,42	1.70	5 (15%)	45,68,68	1.61	8 (17%)

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	EDO	D	305	-	-	1/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	EDO	D	304	-	-	0/1/1/1	-
5	0O2	D	306	3	-	7/29/49/49	0/3/3/3
6	EDO	D	303	-	-	0/1/1/1	-
5	0O2	C	304	3	-	4/29/49/49	0/3/3/3
5	0O2	B	303	3	-	1/29/49/49	0/3/3/3

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	303	0O2	O4'-C1'	6.14	1.49	1.41
5	C	304	0O2	O4'-C1'	5.72	1.49	1.41
5	D	306	0O2	O4'-C1'	5.12	1.48	1.41
5	B	303	0O2	C2'-C1'	-3.76	1.48	1.53
5	D	306	0O2	C2'-C1'	-3.70	1.48	1.53
5	C	304	0O2	C2-N2	3.69	1.41	1.33
5	B	303	0O2	C2-N2	3.39	1.40	1.33
5	D	306	0O2	C2-N2	3.33	1.40	1.33
5	C	304	0O2	C2'-C1'	-3.16	1.49	1.53
5	D	306	0O2	C2'-C3'	-3.07	1.46	1.52
5	B	303	0O2	C2'-C3'	-2.72	1.46	1.52
5	C	304	0O2	C2'-C3'	-2.55	1.47	1.52
5	C	304	0O2	O4'-C4'	2.44	1.50	1.45
5	D	306	0O2	O4'-C4'	2.35	1.50	1.45
5	B	303	0O2	O4'-C4'	2.22	1.50	1.45

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	C	304	0O2	N3-C2-N1	-4.64	121.03	127.22
5	B	303	0O2	N3-C2-N1	-4.58	121.12	127.22
5	D	306	0O2	N3-C2-N1	-4.41	121.34	127.22
5	C	304	0O2	C6-N1-C2	4.05	122.36	115.93
5	B	303	0O2	C6-N1-C2	4.03	122.33	115.93
5	D	306	0O2	C2-N3-C4	3.80	119.70	115.36
5	D	306	0O2	C4-C5-N7	-3.67	105.57	109.40
5	B	303	0O2	C2-N3-C4	3.56	119.42	115.36
5	C	304	0O2	C2-N3-C4	3.41	119.26	115.36
5	C	304	0O2	C5-C6-N1	-3.38	118.81	123.43
5	B	303	0O2	C5-C6-N1	-3.33	118.88	123.43
5	B	303	0O2	C4-C5-N7	-3.30	105.96	109.40
5	D	306	0O2	C6-N1-C2	3.28	121.14	115.93
5	C	304	0O2	C6-C5-C4	-3.17	117.77	120.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	303	0O2	C6-C5-C4	-2.92	118.01	120.80
5	C	304	0O2	N2-C2-N1	2.87	121.71	117.25
5	D	306	0O2	C5-C6-N1	-2.85	119.53	123.43
5	D	306	0O2	C6-C5-C4	-2.82	118.11	120.80
5	C	304	0O2	PB-O3B-PG	-2.73	123.45	132.83
5	C	304	0O2	C4-C5-N7	-2.63	106.66	109.40
5	C	304	0O2	C3'-C2'-C1'	2.46	105.33	99.89
5	C	304	0O2	O4'-C1'-C2'	-2.40	103.42	106.93
5	B	303	0O2	PB-O3B-PG	-2.37	124.69	132.83
5	D	306	0O2	PA-O3A-PB	2.21	140.41	132.83
5	D	306	0O2	O4'-C1'-C2'	-2.13	103.81	106.93
5	D	306	0O2	N2-C2-N1	2.13	120.56	117.25
5	B	303	0O2	O3G-PG-O3B	2.11	111.70	104.64
5	D	306	0O2	O3G-PG-O3B	2.10	111.68	104.64
5	C	304	0O2	O3G-PG-O3B	2.03	111.45	104.64

There are no chirality outliers.

All (13) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	C	304	0O2	PC-O3C-PD-O1D
5	C	304	0O2	PB-O3B-PG-O3G
5	D	306	0O2	PB-O3B-PG-O3G
5	D	306	0O2	PB-O3A-PA-O5'
6	D	305	EDO	O1-C1-C2-O2
5	D	306	0O2	PG-O3B-PB-O2B
5	C	304	0O2	PB-O3B-PG-O2G
5	D	306	0O2	PC-O3C-PD-O1D
5	B	303	0O2	PC-O3C-PD-O1D
5	D	306	0O2	PA-O3A-PB-O1B
5	C	304	0O2	PB-O3B-PG-O1G
5	D	306	0O2	PB-O3B-PG-O2G
5	D	306	0O2	PB-O3A-PA-O2A

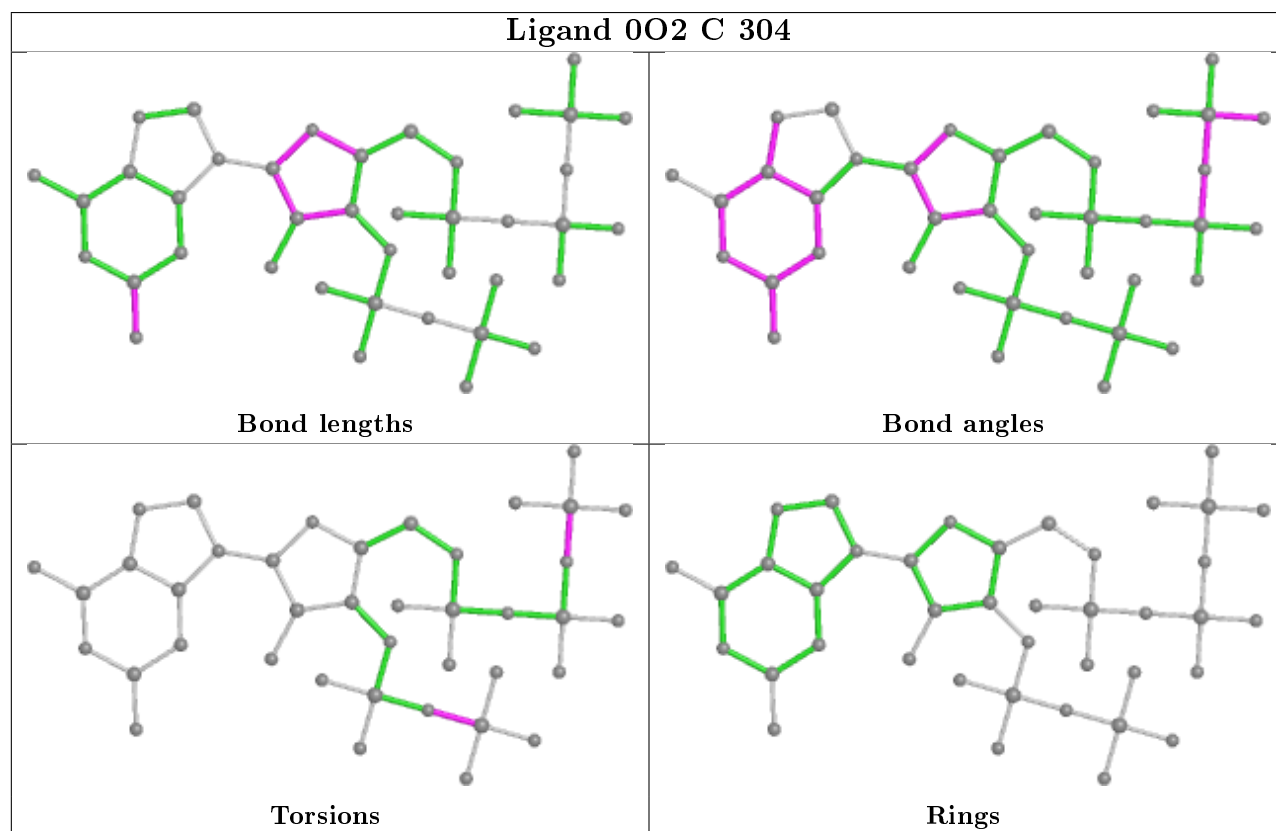
There are no ring outliers.

1 monomer is involved in 3 short contacts:

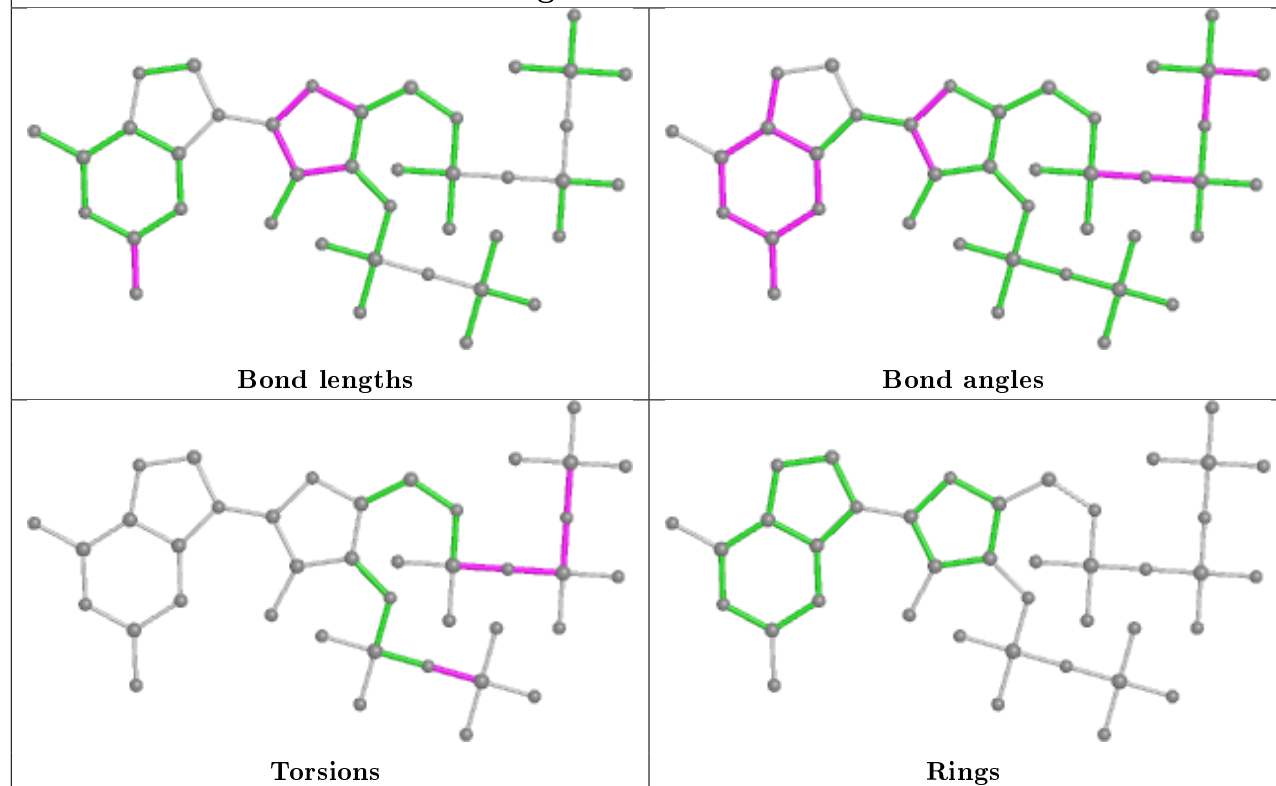
Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	D	305	EDO	3	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths,

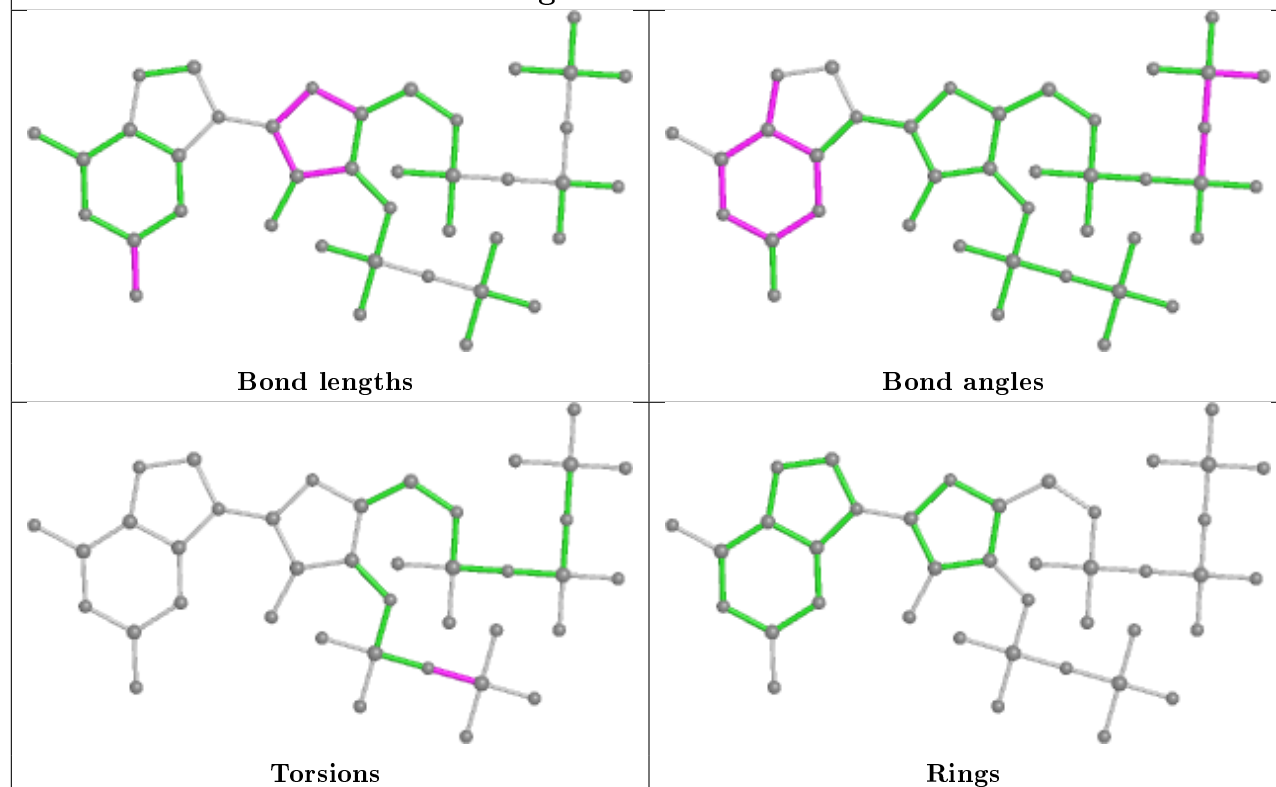
bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 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 0O2 D 306



Ligand 0O2 B 303



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	180/210 (85%)	0.33	14 (7%) 13 12	14, 25, 55, 75	0
1	B	186/210 (88%)	0.33	16 (8%) 10 10	14, 23, 53, 73	0
1	C	185/210 (88%)	0.42	19 (10%) 6 5	14, 24, 63, 89	0
1	D	184/210 (87%)	0.44	16 (8%) 10 9	14, 23, 59, 86	0
All	All	735/840 (87%)	0.38	65 (8%) 10 9	14, 24, 59, 89	0

All (65) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	147	ILE	15.1
1	D	198	TYR	8.2
1	C	151	ILE	8.1
1	A	3	ASN	7.1
1	C	132	LEU	6.8
1	B	131	HIS	6.6
1	C	152	ASN	6.3
1	C	129	LEU	6.2
1	B	132	LEU	6.1
1	B	3	ASN	6.1
1	D	163	LEU	6.0
1	D	3	ASN	5.8
1	D	148	GLN	5.7
1	C	128	SER	5.7
1	A	152	ASN	5.4
1	B	2	ASP	5.1
1	C	3	ASN	5.0
1	C	153	GLU	4.9
1	B	153	GLU	4.7
1	A	131	HIS	4.7
1	C	130	GLU	4.5

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Mol	Chain	Res	Type	RSRZ
1	D	132	LEU	4.5
1	C	131	HIS	4.4
1	A	200	LYS	4.3
1	D	131	HIS	4.3
1	A	132	LEU	4.2
1	B	152	ASN	4.1
1	C	163	LEU	4.0
1	B	47	GLU	3.8
1	B	129	LEU	3.8
1	A	153	GLU	3.7
1	B	154	ALA	3.6
1	C	205	ALA	3.6
1	D	151	ILE	3.5
1	A	201	MET	3.5
1	C	156	LYS	3.4
1	A	199	ARG	3.4
1	D	150	ARG	3.2
1	D	200	LYS	3.2
1	C	162	ASN	3.0
1	B	155	ARG	3.0
1	D	162	ASN	3.0
1	D	155	ARG	3.0
1	D	47	GLU	2.9
1	A	198	TYR	2.6
1	A	156	LYS	2.6
1	A	78	TYR	2.6
1	D	153	GLU	2.5
1	C	158	VAL	2.5
1	B	156	LYS	2.5
1	C	203	LEU	2.4
1	A	4	GLU	2.4
1	C	199	ARG	2.4
1	D	152	ASN	2.4
1	B	201	MET	2.4
1	C	155	ARG	2.3
1	A	44	GLN	2.3
1	B	48	GLY	2.3
1	D	197	LYS	2.3
1	C	204	GLU	2.2
1	B	206	LYS	2.2
1	B	128	SER	2.1
1	B	130	GLU	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	47	GLU	2.1
1	C	157	GLU	2.0

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

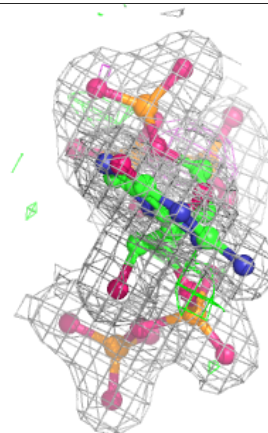
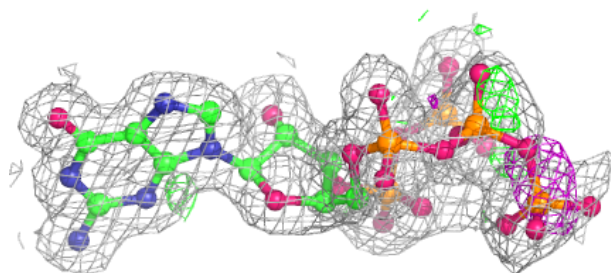
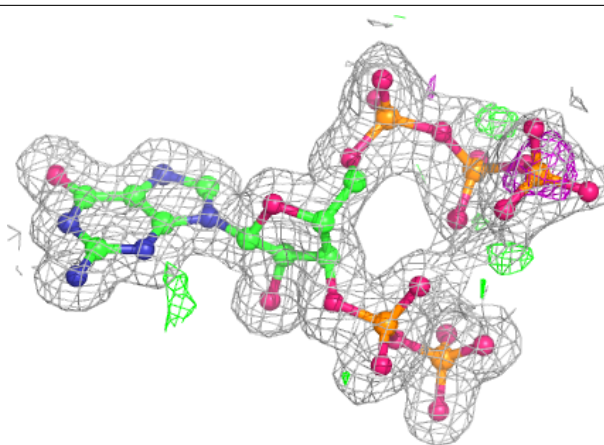
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
6	EDO	D	303	4/4	0.91	0.24	27,27,27,29	0
6	EDO	D	305	4/4	0.93	0.16	19,21,23,23	0
6	EDO	D	304	4/4	0.94	0.09	23,24,24,25	0
5	0O2	B	303	40/40	0.94	0.10	13,22,58,59	0
5	0O2	C	304	40/40	0.95	0.09	14,21,57,59	0
5	0O2	D	306	40/40	0.96	0.10	12,17,55,57	0
3	MG	D	301	1/1	0.96	0.12	23,23,23,23	0
2	SO4	A	301	5/5	0.97	0.07	31,32,34,34	0
3	MG	B	301	1/1	0.98	0.08	21,21,21,21	0
2	SO4	C	303	5/5	0.98	0.06	24,24,25,26	0
3	MG	C	301	1/1	0.99	0.07	22,22,22,22	0
4	K	C	302	1/1	0.99	0.06	15,15,15,15	0
4	K	D	302	1/1	1.00	0.08	15,15,15,15	0
4	K	B	302	1/1	1.00	0.07	14,14,14,14	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.

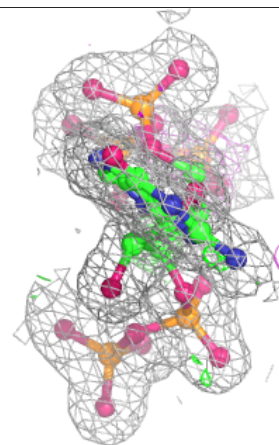
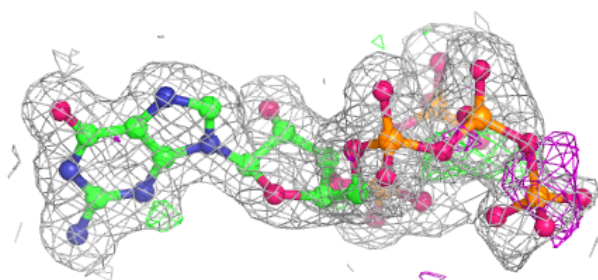
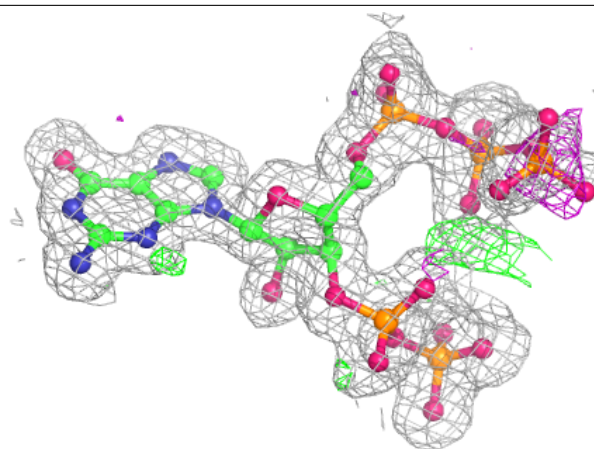
Electron density around 0O2 B 303:

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



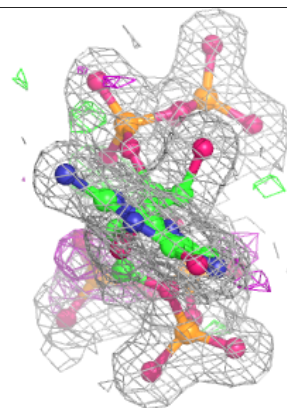
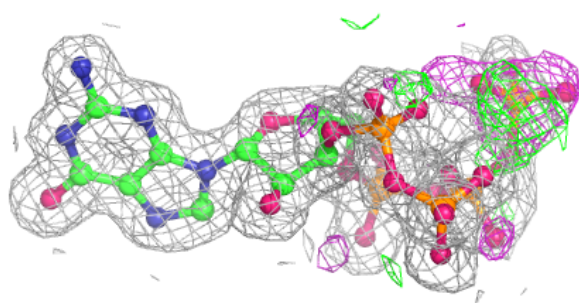
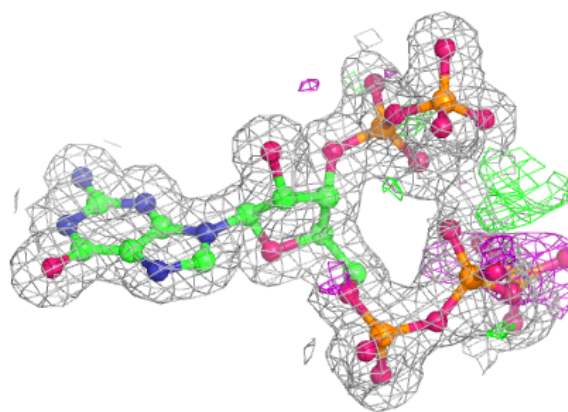
Electron density around 0O2 C 304:

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



Electron density around 0O2 D 306:

$2mF_o-DF_c$ (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.