



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 10:41 AM UTC

PDB ID : 3ZWO / pdb_00003zwo
Title : Crystal structure of ADP ribosyl cyclase complexed with reaction intermediate
Authors : Kotaka, M.; Graeff, R.; Zhang, L.H.; Lee, H.C.; Hao, Q.
Deposited on : 2011-08-02
Resolution : 2.00 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

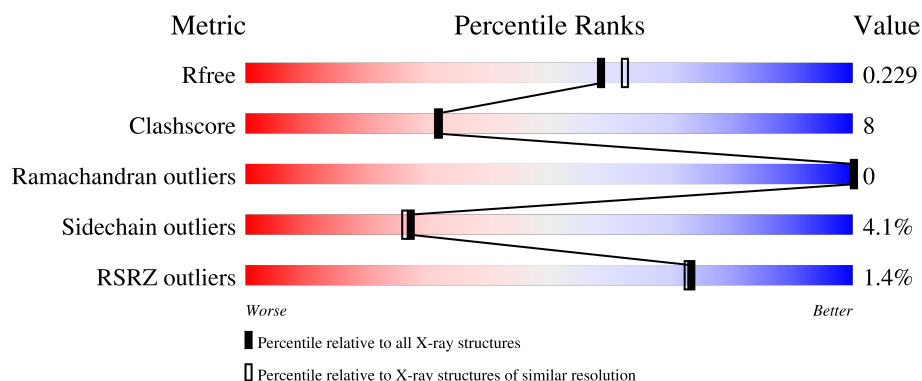
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments 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	259	% 82% 12% . .
1	B	259	2% 84% 12% . .
1	C	259	% 83% 13% . .
1	D	259	% 80% 15% . . .
1	E	259	2% 83% 13% . .

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Mol	Chain	Length	Quality of chain
1	F	259	
1	G	259	
1	H	259	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	G2Q	A	301	-	-	X	-
2	G2Q	B	301	-	-	X	-
2	G2Q	C	301	-	-	X	-
2	G2Q	E	301	-	-	X	-
2	G2Q	G	301	-	-	X	-

2 Entry composition

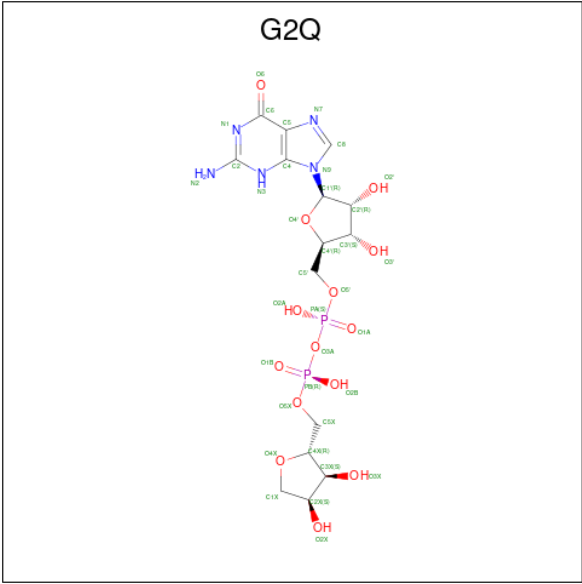
There are 4 unique types of molecules in this entry. The entry contains 17700 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 ADP-RIBOSYL CYCLASE.

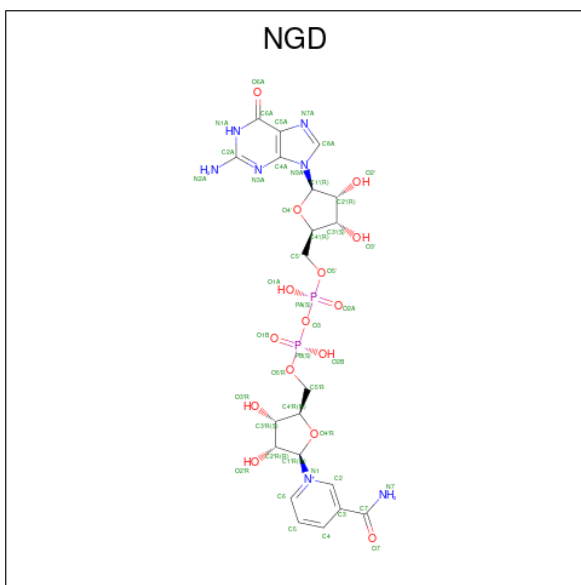
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	252	Total	C	N	O	S	0	0	0
			2017	1291	343	369	14			
1	B	251	Total	C	N	O	S	0	0	0
			2012	1288	342	368	14			
1	C	252	Total	C	N	O	S	0	0	0
			2017	1291	343	369	14			
1	D	251	Total	C	N	O	S	0	0	0
			2012	1288	342	368	14			
1	E	251	Total	C	N	O	S	0	0	0
			2012	1288	342	368	14			
1	F	251	Total	C	N	O	S	0	0	0
			2012	1288	342	368	14			
1	G	252	Total	C	N	O	S	0	0	0
			2017	1291	343	369	14			
1	H	251	Total	C	N	O	S	0	0	0
			2012	1288	342	368	14			

- Molecule 2 is GUANOSINE DIPHOSPHATE RIBOSE (CCD ID: G2Q) (formula: $C_{15}H_{23}N_5O_{14}P_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			36	15	5	14	2		
2	B	1	Total	C	N	O	P	0	0
			36	15	5	14	2		
2	C	1	Total	C	N	O	P	0	0
			36	15	5	14	2		
2	E	1	Total	C	N	O	P	0	0
			36	15	5	14	2		
2	G	1	Total	C	N	O	P	0	0
			36	15	5	14	2		
2	H	1	Total	C	N	O	P	0	0
			36	15	5	14	2		

- Molecule 3 is 3-(AMINOCARBONYL)-1-[(2R,3R,4S,5R)-5-({[(S)-{(S)-[(2R,3S,4R,5R)-5-(2-AMINO-6-OXO-1,6-DIHYDRO-9H-PURIN-9-YL)-3,4-DIHYDROXYTETRAHYDROFURAN-2-YL]METHOXY}(HYDROXY)PHOSPHORYL]OXY}(HYDROXY)PHOSPHORYL]OXY}METHYL)-3,4-DIHYDROXYTETRAHYDROFURAN-2-YL]PYRIDINIUM (CCD ID: NGD) (formula: C₂₁H₂₈N₇O₁₅P₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	F	1	Total 45	C 21	N 7	O 15	P 2	0	0

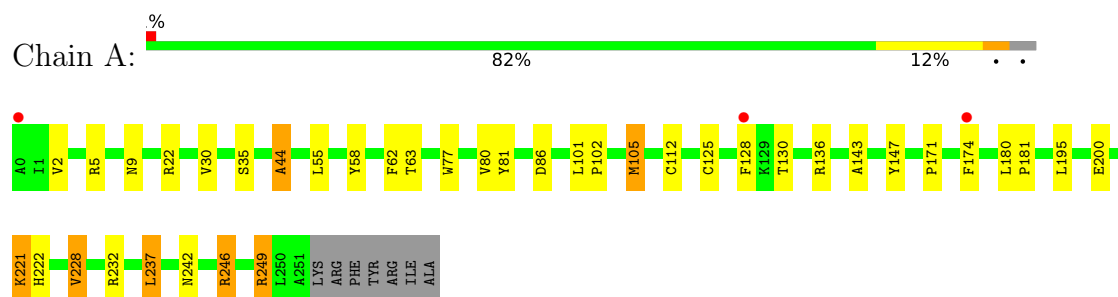
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	204	Total O 204 204	0	0
4	B	194	Total O 194 194	0	0
4	C	208	Total O 208 208	0	0
4	D	204	Total O 204 204	0	0
4	E	182	Total O 182 182	0	0
4	F	138	Total O 138 138	0	0
4	G	100	Total O 100 100	0	0
4	H	98	Total O 98 98	0	0

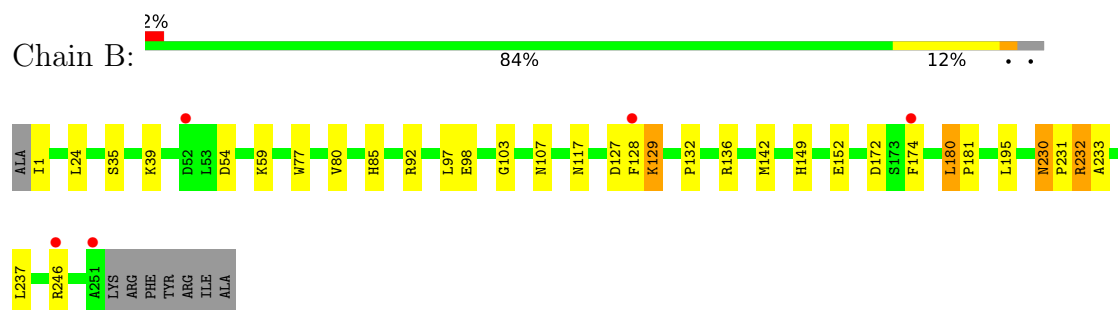
3 Residue-property plots [i](#)

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.

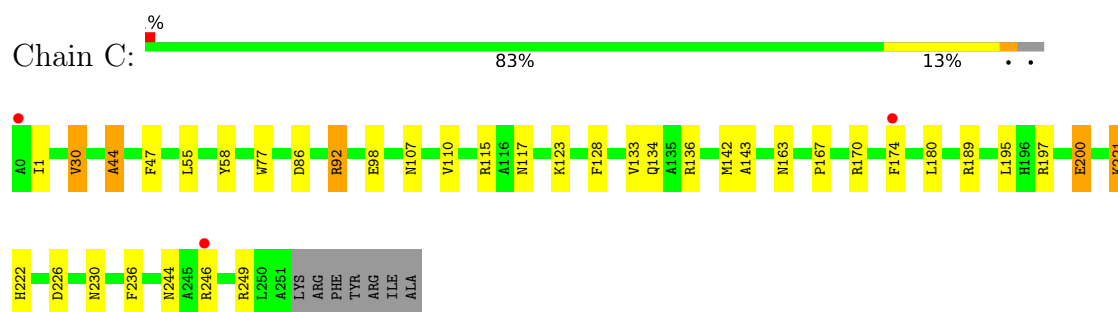
• Molecule 1: ADP-RIBOSYL CYCLASE



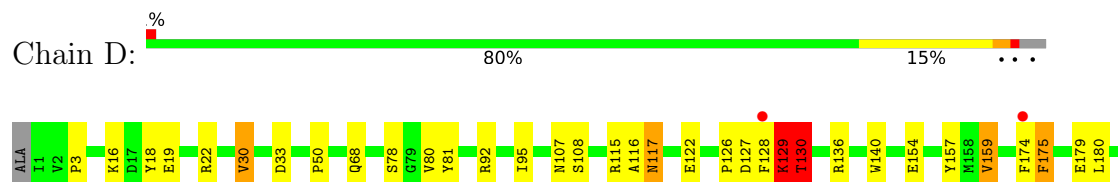
• Molecule 1: ADP-RIBOSYL CYCLASE



• Molecule 1: ADP-RIBOSYL CYCLASE

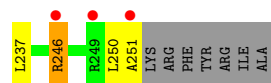
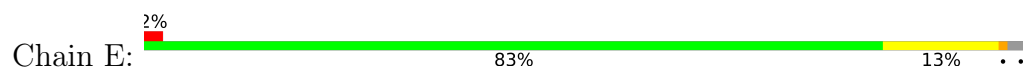


• Molecule 1: ADP-RIBOSYL CYCLASE

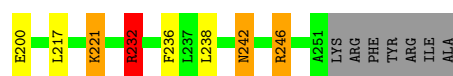
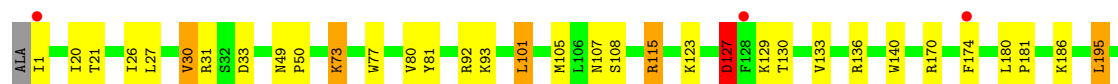
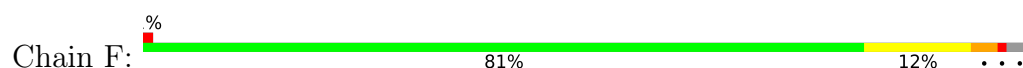




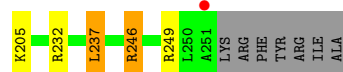
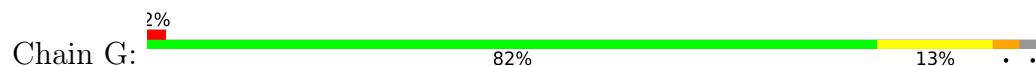
• Molecule 1: ADP-RIBOSYL CYCLASE



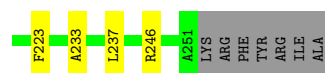
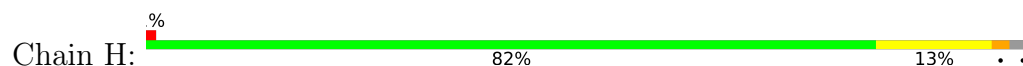
• Molecule 1: ADP-RIBOSYL CYCLASE



• Molecule 1: ADP-RIBOSYL CYCLASE



• Molecule 1: ADP-RIBOSYL CYCLASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	60.57Å 76.42Å 139.71Å 87.80° 89.09° 89.32°	Depositor
Resolution (Å)	30.00 – 2.00 30.00 – 2.00	Depositor EDS
% Data completeness (in resolution range)	97.1 (30.00-2.00) 83.4 (30.00-2.00)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.53 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.177 , 0.228 (Not available) , 0.229	Depositor DCC
R_{free} test set	7075 reflections (4.95%)	wwPDB-VP
Wilson B-factor (Å ²)	31.0	Xtriage
Anisotropy	0.107	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 44.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.059 for h,-k,-l 0.015 for -h,k,-l 0.004 for -h,-k,l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	17700	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.08% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: NGD, G2Q

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	1.43	6/2069 (0.3%)	1.25	6/2800 (0.2%)
1	B	1.40	7/2064 (0.3%)	1.23	6/2793 (0.2%)
1	C	1.45	7/2069 (0.3%)	1.24	6/2800 (0.2%)
1	D	1.35	3/2064 (0.1%)	1.21	6/2793 (0.2%)
1	E	1.39	3/2064 (0.1%)	1.22	5/2793 (0.2%)
1	F	1.27	3/2064 (0.1%)	1.19	7/2793 (0.3%)
1	G	1.12	1/2069 (0.0%)	1.10	6/2800 (0.2%)
1	H	1.10	3/2064 (0.1%)	1.07	3/2793 (0.1%)
All	All	1.32	33/16527 (0.2%)	1.19	45/22365 (0.2%)

All (33) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	44	ALA	CA-CB	8.20	1.66	1.53
1	C	44	ALA	CA-CB	7.53	1.65	1.53
1	B	233	ALA	CA-CB	7.18	1.65	1.53
1	C	230	ASN	CA-CB	6.80	1.58	1.52
1	A	105	MET	N-CA	6.49	1.54	1.46
1	C	143	ALA	CA-CB	6.09	1.62	1.53
1	F	30	VAL	CA-C	5.94	1.59	1.53
1	B	132	PRO	C-O	5.92	1.30	1.23
1	B	80	VAL	CA-CB	5.68	1.62	1.54
1	A	63	THR	CA-CB	5.68	1.62	1.53
1	C	133	VAL	CA-CB	5.68	1.61	1.54
1	H	117	ASN	CA-C	5.61	1.59	1.52
1	D	230	ASN	N-CA	5.53	1.50	1.46
1	C	107	ASN	N-CA	5.46	1.53	1.46
1	C	110	VAL	CB-CG2	5.44	1.70	1.52
1	F	21	THR	CA-CB	5.35	1.61	1.54
1	A	147	TYR	C-O	5.35	1.30	1.24
1	E	169	TYR	N-CA	5.33	1.52	1.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	97	LEU	CA-CB	5.29	1.61	1.53
1	C	167	PRO	CA-C	5.29	1.58	1.52
1	D	95	ILE	N-CA	5.27	1.52	1.46
1	B	152	GLU	N-CA	5.25	1.52	1.45
1	D	180	LEU	CA-C	5.25	1.58	1.52
1	H	233	ALA	CA-C	5.24	1.59	1.52
1	E	230	ASN	CA-CB	5.21	1.57	1.52
1	A	2	VAL	N-CA	5.20	1.51	1.46
1	B	230	ASN	CA-CB	5.18	1.57	1.52
1	E	49	ASN	CB-CG	5.18	1.65	1.52
1	A	130	THR	CA-CB	5.17	1.62	1.53
1	H	81	TYR	N-CA	5.15	1.52	1.46
1	B	180	LEU	C-O	5.10	1.29	1.24
1	G	43	LYS	N-CA	5.09	1.52	1.46
1	F	27	LEU	C-N	5.08	1.39	1.33

All (45) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	24	LEU	CA-C-N	-8.88	108.56	122.67
1	B	24	LEU	C-N-CA	-8.88	108.56	122.67
1	D	129	LYS	N-CA-C	-8.08	102.89	114.12
1	A	2	VAL	N-CA-C	7.79	116.50	107.77
1	E	24	LEU	CA-C-N	-7.36	110.87	122.73
1	E	24	LEU	C-N-CA	-7.36	110.87	122.73
1	G	43	LYS	CB-CA-C	-7.32	98.23	110.68
1	H	127	ASP	N-CA-C	-7.20	99.24	109.96
1	B	230	ASN	CA-C-N	-6.66	112.80	119.92
1	B	230	ASN	C-N-CA	-6.66	112.80	119.92
1	E	203	ILE	N-CA-C	-6.38	106.32	111.81
1	F	127	ASP	N-CA-C	-6.36	100.24	109.59
1	C	30	VAL	N-CA-CB	-6.36	99.90	112.24
1	A	125	CYS	CA-C-N	-6.35	114.09	120.31
1	A	125	CYS	C-N-CA	-6.35	114.09	120.31
1	D	30	VAL	N-CA-CB	-6.21	100.00	111.91
1	F	30	VAL	N-CA-CB	-6.05	101.34	111.38
1	C	180	LEU	CA-C-N	-5.85	113.66	119.56
1	C	180	LEU	C-N-CA	-5.85	113.66	119.56
1	E	232	ARG	N-CA-C	5.74	118.00	111.11
1	F	30	VAL	CB-CA-C	5.74	117.90	111.59
1	G	69	LEU	CA-C-N	-5.64	114.16	120.14
1	G	69	LEU	C-N-CA	-5.64	114.16	120.14

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	175	PHE	N-CA-C	5.61	117.07	111.07
1	E	172	ASP	N-CA-C	5.60	120.11	113.16
1	B	172	ASP	N-CA-C	5.47	119.82	113.20
1	G	27	LEU	CA-C-N	-5.45	115.13	120.52
1	G	27	LEU	C-N-CA	-5.45	115.13	120.52
1	D	19	GLU	N-CA-C	-5.44	104.58	111.11
1	H	170	ARG	CA-C-N	-5.43	114.07	119.56
1	H	170	ARG	C-N-CA	-5.43	114.07	119.56
1	F	232	ARG	N-CA-C	5.42	117.89	111.33
1	C	163	ASN	CA-C-N	-5.40	114.26	119.87
1	C	163	ASN	C-N-CA	-5.40	114.26	119.87
1	B	85	HIS	N-CA-C	5.32	117.16	111.36
1	C	170	ARG	CB-CA-C	-5.31	101.21	109.55
1	D	232	ARG	N-CA-C	5.25	117.41	111.11
1	F	101	LEU	CA-C-N	-5.23	113.15	118.85
1	F	101	LEU	C-N-CA	-5.23	113.15	118.85
1	A	112	CYS	CA-C-N	5.22	125.30	120.34
1	A	112	CYS	C-N-CA	5.22	125.30	120.34
1	D	130	THR	N-CA-C	5.22	117.87	111.82
1	F	242	ASN	N-CA-CB	-5.12	104.92	110.71
1	G	133	VAL	N-CA-C	5.04	115.77	110.62
1	A	22	ARG	N-CA-C	5.02	117.99	109.76

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	2017	0	1973	25	0
1	B	2012	0	1968	16	0
1	C	2017	0	1973	25	0
1	D	2012	0	1968	39	1
1	E	2012	0	1968	28	0
1	F	2012	0	1968	35	1
1	G	2017	0	1973	28	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	H	2012	0	1968	31	0
2	A	36	0	21	10	0
2	B	36	0	21	12	0
2	C	36	0	21	13	0
2	E	36	0	21	11	0
2	G	36	0	21	9	0
2	H	36	0	21	5	0
3	F	45	0	26	5	0
4	A	204	0	0	9	0
4	B	194	0	0	10	0
4	C	208	0	0	12	0
4	D	204	0	0	12	1
4	E	182	0	0	14	0
4	F	138	0	0	8	1
4	G	100	0	0	6	0
4	H	98	0	0	3	0
All	All	17700	0	15911	262	2

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 8.

All (262) 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)
2:C:301:G2Q:H1XA	4:C:2106:HOH:O	1.25	1.33
1:A:86:ASP:HB3	4:A:2108:HOH:O	1.26	1.30
1:C:86:ASP:HB3	4:C:2116:HOH:O	1.24	1.25
1:D:200:GLU:HG2	4:D:2168:HOH:O	1.41	1.19
1:F:49:ASN:HB2	1:F:115:ARG:HH22	1.05	1.13
1:G:127:ASP:O	1:G:130:THR:HG22	1.49	1.13
1:D:174:PHE:HZ	4:D:2055:HOH:O	1.34	1.10
1:F:200:GLU:HG2	4:F:2120:HOH:O	1.50	1.10
2:C:301:G2Q:C1X	4:C:2106:HOH:O	1.86	1.09
1:F:246:ARG:HH11	1:F:246:ARG:HG2	1.14	1.06
1:A:232:ARG:NH1	4:A:2187:HOH:O	1.85	1.06
1:E:92:ARG:HD2	4:E:2101:HOH:O	1.58	1.03
1:F:246:ARG:HH11	1:F:246:ARG:CG	1.71	1.02
1:D:68:GLN:CG	4:D:2081:HOH:O	2.13	0.95
1:F:49:ASN:HB2	1:F:115:ARG:NH2	1.80	0.95
1:D:68:GLN:HG3	4:D:2081:HOH:O	1.67	0.94
1:G:174:PHE:CZ	4:G:2100:HOH:O	2.19	0.94

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:55:LEU:HD22	4:C:2080:HOH:O	1.71	0.90
1:D:117:ASN:H	1:D:117:ASN:HD22	1.14	0.89
1:C:55:LEU:CD2	4:C:2080:HOH:O	2.21	0.88
1:H:132:PRO:HB2	1:H:134:GLN:HE22	1.39	0.88
1:H:132:PRO:HB2	1:H:134:GLN:NE2	1.89	0.87
1:G:174:PHE:HZ	4:G:2100:HOH:O	1.52	0.87
1:H:174:PHE:CE2	2:H:301:G2Q:O2'	2.28	0.86
1:E:174:PHE:CE1	4:E:2129:HOH:O	2.28	0.86
1:F:93:LYS:HE2	4:F:2079:HOH:O	1.76	0.86
1:E:77:TRP:O	2:E:301:G2Q:H5XA	1.77	0.84
1:C:142:MET:HE1	4:C:2156:HOH:O	1.77	0.83
2:C:301:G2Q:C2X	4:C:2106:HOH:O	2.19	0.82
1:C:92:ARG:HG3	1:C:92:ARG:HH11	1.44	0.81
1:H:117:ASN:HD22	1:H:118:PRO:HA	1.45	0.81
1:F:246:ARG:CG	1:F:246:ARG:NH1	2.38	0.79
1:F:26:ILE:HG13	4:F:2026:HOH:O	1.82	0.78
1:G:127:ASP:O	1:G:130:THR:CG2	2.30	0.77
1:B:77:TRP:O	2:B:301:G2Q:H5XA	1.83	0.77
1:E:222:HIS:ND1	4:E:2171:HOH:O	2.18	0.76
1:E:233:ALA:O	4:E:2177:HOH:O	2.03	0.76
2:A:301:G2Q:O4X	2:A:301:G2Q:O3'	2.02	0.75
1:B:246:ARG:HG3	4:B:2194:HOH:O	1.84	0.75
1:E:49:ASN:HB2	4:E:2059:HOH:O	1.85	0.75
1:H:54:ASP:HB2	4:H:2044:HOH:O	1.86	0.75
1:B:127:ASP:OD1	1:B:129:LYS:HG2	1.86	0.74
1:H:137:GLU:HG2	1:H:174:PHE:CE1	2.23	0.74
2:B:301:G2Q:O3'	2:B:301:G2Q:H1XA	1.90	0.72
1:H:132:PRO:CB	1:H:134:GLN:HE22	2.02	0.72
1:B:92:ARG:NH1	4:B:2109:HOH:O	2.22	0.71
4:B:2193:HOH:O	1:G:232:ARG:NH2	2.23	0.71
2:E:301:G2Q:O3'	2:E:301:G2Q:O4X	2.07	0.71
1:C:77:TRP:O	2:C:301:G2Q:H5XA	1.90	0.71
1:A:77:TRP:O	2:A:301:G2Q:H5XA	1.89	0.70
2:C:301:G2Q:H2X	4:C:2106:HOH:O	1.88	0.70
1:E:47:PHE:HE1	1:E:117:ASN:HD22	1.37	0.69
1:H:126:PRO:HB2	1:H:130:THR:HG23	1.76	0.68
1:F:246:ARG:NH2	4:F:2137:HOH:O	2.14	0.68
1:E:174:PHE:CD1	4:E:2129:HOH:O	2.46	0.68
1:C:244:ASN:OD1	1:F:232:ARG:NH2	2.26	0.68
1:D:117:ASN:H	1:D:117:ASN:ND2	1.89	0.67
2:E:301:G2Q:O3'	2:E:301:G2Q:C1X	2.43	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:301:G2Q:O3'	2:B:301:G2Q:C1X	2.43	0.66
1:E:47:PHE:HE1	1:E:117:ASN:ND2	1.94	0.65
1:D:117:ASN:ND2	4:D:2123:HOH:O	2.29	0.65
1:F:246:ARG:HG2	1:F:246:ARG:NH1	1.97	0.65
1:H:137:GLU:HG2	1:H:174:PHE:CZ	2.31	0.65
1:A:86:ASP:CB	4:A:2108:HOH:O	2.06	0.65
1:D:154:GLU:OE2	4:D:2136:HOH:O	2.14	0.65
1:H:127:ASP:O	1:H:130:THR:HG22	1.96	0.65
1:G:25:ASP:HB2	4:G:2022:HOH:O	1.95	0.65
2:C:301:G2Q:H2'	2:C:301:G2Q:HN3	1.62	0.64
1:C:92:ARG:HG3	1:C:92:ARG:NH1	2.06	0.64
1:E:6:GLU:OE1	4:E:2011:HOH:O	2.15	0.64
1:D:128:PHE:O	1:D:136:ARG:HD3	1.98	0.64
1:G:79:GLY:N	2:G:301:G2Q:O2A	2.31	0.63
1:B:174:PHE:CD2	2:B:301:G2Q:O2'	2.51	0.63
1:F:242:ASN:ND2	4:F:2134:HOH:O	2.21	0.63
1:F:246:ARG:NH1	4:F:2137:HOH:O	2.22	0.63
1:H:127:ASP:O	1:H:130:THR:CG2	2.47	0.63
2:C:301:G2Q:O3'	2:C:301:G2Q:H5X	1.98	0.63
1:D:127:ASP:O	1:D:130:THR:CG2	2.47	0.63
1:F:174:PHE:CD2	3:F:5573:NGD:H6	2.35	0.61
1:D:127:ASP:O	1:D:129:LYS:N	2.31	0.61
1:F:127:ASP:O	1:F:130:THR:HG23	2.00	0.61
1:H:174:PHE:CZ	2:H:301:G2Q:O2'	2.51	0.61
1:E:186:LYS:HD2	4:E:2132:HOH:O	2.01	0.61
1:F:77:TRP:O	3:F:5573:NGD:H5'1	2.01	0.60
1:F:200:GLU:CG	4:F:2120:HOH:O	2.24	0.60
1:C:92:ARG:HH11	1:C:92:ARG:CG	2.13	0.59
1:E:49:ASN:CG	1:E:50:PRO:HD2	2.27	0.59
1:G:31:ARG:HD3	4:G:2027:HOH:O	2.02	0.59
1:D:78:SER:HB2	1:D:159:VAL:HG12	1.83	0.59
1:A:5:ARG:HD2	4:A:2011:HOH:O	2.02	0.59
1:G:174:PHE:CD2	2:G:301:G2Q:H2'	2.38	0.59
1:C:115:ARG:NH1	4:C:2072:HOH:O	2.36	0.58
1:F:174:PHE:HB2	3:F:5573:NGD:PB	2.43	0.58
1:G:80:VAL:HG21	1:G:158:MET:HG2	1.85	0.58
1:H:68:GLN:HG3	4:H:2049:HOH:O	2.04	0.57
2:C:301:G2Q:H2'	2:C:301:G2Q:N3	2.18	0.57
2:G:301:G2Q:H1XA	2:G:301:G2Q:N7	2.20	0.56
2:A:301:G2Q:O3'	2:A:301:G2Q:C5X	2.53	0.56
1:C:244:ASN:HB2	4:C:2202:HOH:O	2.05	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:301:G2Q:O3'	2:E:301:G2Q:H1XA	2.06	0.56
1:G:103:GLY:O	1:G:107:ASN:HB2	2.06	0.56
1:A:174:PHE:CD2	2:A:301:G2Q:O2'	2.57	0.55
1:A:237:LEU:HD13	1:D:240:SER:CB	2.36	0.55
1:C:55:LEU:HD22	1:C:55:LEU:H	1.71	0.55
1:D:127:ASP:O	1:D:130:THR:HG22	2.07	0.55
2:C:301:G2Q:O3'	2:C:301:G2Q:C5X	2.54	0.55
1:A:237:LEU:HD13	1:D:240:SER:HB2	1.88	0.55
1:E:92:ARG:CD	4:E:2101:HOH:O	2.33	0.54
1:G:246:ARG:O	1:G:249:ARG:HG3	2.08	0.54
1:C:98:GLU:OE2	2:C:301:G2Q:N2	2.41	0.54
1:E:174:PHE:CD1	2:E:301:G2Q:H4'	2.43	0.54
1:D:117:ASN:CG	4:D:2124:HOH:O	2.51	0.54
1:G:0:ALA:HB1	1:G:124:VAL:HG22	1.90	0.54
1:D:127:ASP:O	1:D:130:THR:HG23	2.07	0.54
1:B:180:LEU:HB3	1:B:181:PRO:HD3	1.90	0.53
1:F:217:LEU:O	1:F:221:LYS:HG2	2.07	0.53
1:B:128:PHE:HD2	1:B:136:ARG:HA	1.74	0.53
1:C:197:ARG:HB2	1:C:200:GLU:HG3	1.91	0.53
2:A:301:G2Q:O3'	2:A:301:G2Q:C1X	2.57	0.53
1:D:117:ASN:HD22	1:D:117:ASN:N	1.95	0.53
2:B:301:G2Q:HN3	2:B:301:G2Q:C2'	2.22	0.52
2:G:301:G2Q:H1XA	2:G:301:G2Q:C8	2.40	0.52
1:D:127:ASP:C	1:D:129:LYS:H	2.17	0.52
1:D:200:GLU:CG	4:D:2168:HOH:O	2.22	0.52
1:A:246:ARG:O	1:A:249:ARG:HG3	2.10	0.52
1:F:49:ASN:OD1	1:F:50:PRO:HD2	2.10	0.52
1:G:136:ARG:HG2	1:G:137:GLU:HG3	1.92	0.52
1:E:174:PHE:CZ	2:E:301:G2Q:H1'	2.46	0.51
1:H:101:LEU:HG	1:H:105:MET:HE2	1.92	0.51
1:G:44:ALA:O	1:G:48:LYS:NZ	2.43	0.51
1:B:232:ARG:NH1	4:B:2188:HOH:O	2.44	0.50
1:G:86:ASP:HB3	1:G:237:LEU:HD21	1.94	0.50
2:B:301:G2Q:HN3	2:B:301:G2Q:C3'	2.25	0.50
4:B:2101:HOH:O	1:G:92:ARG:CD	2.59	0.50
1:G:174:PHE:CG	2:G:301:G2Q:H2'	2.47	0.50
1:E:101:LEU:HG	1:E:105:MET:HE2	1.94	0.50
2:B:301:G2Q:HN3	2:B:301:G2Q:H2'	1.77	0.49
1:B:149:HIS:ND1	4:B:2142:HOH:O	2.34	0.49
4:B:2101:HOH:O	1:G:92:ARG:HD3	2.12	0.49
1:H:218:VAL:HG13	1:H:223:PHE:HB2	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:301:G2Q:O3'	2:B:301:G2Q:O4X	2.30	0.49
1:E:246:ARG:HG3	4:E:2182:HOH:O	2.12	0.49
2:H:301:G2Q:H8	2:H:301:G2Q:H1XA	1.95	0.49
1:A:171:PRO:HG3	1:E:129:LYS:HD3	1.95	0.48
1:C:47:PHE:HE1	1:C:117:ASN:HD22	1.59	0.48
1:H:35:SER:HB3	4:H:2004:HOH:O	2.13	0.48
1:H:47:PHE:HA	1:H:114:GLN:O	2.13	0.48
2:H:301:G2Q:H1XA	2:H:301:G2Q:C8	2.44	0.48
2:B:301:G2Q:HN3	2:B:301:G2Q:H3'	1.78	0.47
1:G:166:VAL:HG13	1:G:170:ARG:HD3	1.95	0.47
1:H:128:PHE:O	1:H:131:CYS:N	2.42	0.47
1:A:128:PHE:HD2	1:A:136:ARG:HA	1.79	0.47
1:F:246:ARG:NH1	1:F:246:ARG:HG3	2.23	0.47
1:G:132:PRO:HB2	1:G:134:GLN:NE2	2.30	0.47
1:C:55:LEU:CD2	1:C:134:GLN:HB2	2.45	0.47
1:D:175:PHE:HA	1:D:179:GLU:HB2	1.97	0.47
1:F:195:LEU:HD11	1:F:238:LEU:HD21	1.97	0.47
2:B:301:G2Q:H2'	2:B:301:G2Q:N3	2.30	0.47
1:C:174:PHE:HZ	4:C:2153:HOH:O	1.97	0.47
1:E:49:ASN:CB	4:E:2059:HOH:O	2.54	0.47
1:F:101:LEU:HG	1:F:105:MET:HE2	1.96	0.47
2:A:301:G2Q:HN3	2:A:301:G2Q:C2'	2.28	0.47
1:C:55:LEU:HD23	4:C:2080:HOH:O	2.04	0.47
1:C:128:PHE:HD2	1:C:136:ARG:HA	1.80	0.46
2:A:301:G2Q:N3	2:A:301:G2Q:H2'	2.29	0.46
1:B:59:LYS:HA	1:B:142:MET:HE2	1.96	0.46
2:A:301:G2Q:HN3	2:A:301:G2Q:H2'	1.80	0.46
2:G:301:G2Q:O5X	2:G:301:G2Q:H5'A	2.16	0.46
1:F:140:TRP:C	1:F:174:PHE:HE2	2.24	0.46
1:B:98:GLU:OE2	2:B:301:G2Q:N2	2.49	0.46
1:D:129:LYS:HE2	1:D:129:LYS:HB2	1.37	0.46
1:D:244:ASN:O	1:D:245:ALA:C	2.58	0.46
2:E:301:G2Q:HN3	2:E:301:G2Q:C2'	2.29	0.46
1:A:44:ALA:HB3	1:A:58:TYR:CE1	2.51	0.46
1:C:189:ARG:HD3	1:C:226:ASP:OD2	2.16	0.46
1:H:98:GLU:OE2	2:H:301:G2Q:O6	2.34	0.45
1:D:50:PRO:HG3	4:D:2121:HOH:O	2.16	0.45
1:D:222:HIS:ND1	4:D:2179:HOH:O	2.36	0.45
1:A:55:LEU:HB2	4:A:2075:HOH:O	2.16	0.45
1:D:116:ALA:O	1:D:117:ASN:C	2.60	0.45
1:A:62:PHE:CZ	1:A:143:ALA:HB2	2.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:174:PHE:HB2	3:F:5573:NGD:O5'R	2.17	0.45
1:B:230:ASN:O	1:B:231:PRO:C	2.59	0.45
1:A:80:VAL:O	1:A:81:TYR:C	2.60	0.44
2:G:301:G2Q:H3'	2:G:301:G2Q:O2B	2.16	0.44
1:H:127:ASP:CG	1:H:129:LYS:HB3	2.42	0.44
1:C:246:ARG:O	1:C:249:ARG:HG3	2.17	0.44
1:H:126:PRO:HB2	1:H:130:THR:CG2	2.44	0.44
2:C:301:G2Q:HN3	2:C:301:G2Q:C2'	2.28	0.44
1:D:68:GLN:HG2	4:D:2081:HOH:O	1.98	0.44
1:A:242:ASN:ND2	4:A:2196:HOH:O	2.40	0.44
1:D:107:ASN:O	1:D:108:SER:HB2	2.18	0.44
1:H:127:ASP:OD2	1:H:129:LYS:HB3	2.18	0.44
1:B:142:MET:HE3	4:B:2072:HOH:O	2.17	0.44
2:E:301:G2Q:HN3	2:E:301:G2Q:H3'	1.82	0.44
1:D:157:TYR:CZ	1:D:159:VAL:CG1	3.01	0.44
1:E:55:LEU:HB2	4:E:2067:HOH:O	2.17	0.43
1:G:39:LYS:O	1:G:43:LYS:HG3	2.18	0.43
1:H:117:ASN:ND2	1:H:118:PRO:HA	2.23	0.43
1:E:77:TRP:NE1	2:E:301:G2Q:H2X	2.33	0.43
1:F:107:ASN:O	1:F:108:SER:HB2	2.19	0.43
1:F:140:TRP:C	1:F:174:PHE:CE2	2.96	0.43
1:E:180:LEU:HB3	1:E:181:PRO:HD3	2.00	0.43
1:F:73:LYS:HA	1:F:73:LYS:HD3	1.82	0.43
1:G:174:PHE:CE2	4:G:2100:HOH:O	2.60	0.43
1:H:86:ASP:HB3	1:H:237:LEU:HD11	1.99	0.43
1:H:101:LEU:O	1:H:105:MET:HG3	2.19	0.43
1:F:170:ARG:HH22	3:F:5573:NGD:H2A2	1.66	0.43
2:E:301:G2Q:HN3	2:E:301:G2Q:H2'	1.83	0.43
1:C:174:PHE:CD2	2:C:301:G2Q:O2'	2.71	0.43
1:A:9:ASN:HB3	1:D:16:LYS:HG2	1.99	0.43
1:C:236:PHE:HB3	1:F:236:PHE:O	2.19	0.43
1:A:174:PHE:CE2	4:A:2147:HOH:O	2.69	0.43
1:B:54:ASP:HB2	4:B:2065:HOH:O	2.18	0.43
1:G:132:PRO:HB2	1:G:134:GLN:HE22	1.84	0.43
1:D:200:GLU:CD	4:D:2168:HOH:O	2.59	0.42
1:E:80:VAL:O	1:E:81:TYR:C	2.62	0.42
2:G:301:G2Q:H3'	2:G:301:G2Q:PB	2.59	0.42
1:E:174:PHE:HZ	4:E:2126:HOH:O	1.97	0.42
1:B:117:ASN:HB2	4:B:2124:HOH:O	2.18	0.42
1:C:221:LYS:O	1:C:222:HIS:HB2	2.20	0.42
1:F:49:ASN:CB	1:F:115:ARG:NH2	2.67	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:126:PRO:O	1:D:128:PHE:CD1	2.72	0.42
1:D:140:TRP:HB3	1:D:174:PHE:HE2	1.84	0.42
1:H:15:CYS:HB2	1:H:105:MET:SD	2.60	0.42
2:E:301:G2Q:H2'	2:E:301:G2Q:N3	2.35	0.42
1:F:246:ARG:CZ	4:F:2137:HOH:O	2.51	0.42
1:G:71:LYS:O	1:G:72:ASN:HB2	2.20	0.42
1:D:128:PHE:HB3	1:D:136:ARG:CG	2.50	0.42
1:A:101:LEU:HG	1:A:105:MET:HE2	2.01	0.42
2:A:301:G2Q:HN3	2:A:301:G2Q:H3'	1.84	0.42
2:B:301:G2Q:C2'	2:B:301:G2Q:N3	2.81	0.42
1:F:180:LEU:HB3	1:F:181:PRO:HD3	2.01	0.42
1:H:127:ASP:O	1:H:130:THR:HG23	2.17	0.42
1:E:250:LEU:O	1:E:251:ALA:C	2.63	0.41
1:A:44:ALA:CB	1:A:58:TYR:CD1	3.03	0.41
1:D:18:TYR:HA	1:D:22:ARG:CG	2.51	0.41
1:G:191:LYS:HE3	1:G:191:LYS:HB3	1.76	0.41
1:A:232:ARG:NH2	4:A:2188:HOH:O	2.53	0.41
1:B:103:GLY:O	1:B:107:ASN:HB2	2.21	0.41
2:C:301:G2Q:H1XA	2:C:301:G2Q:O3'	2.21	0.41
1:H:112:CYS:HB3	1:H:128:PHE:HZ	1.84	0.41
1:F:133:VAL:HB	1:F:136:ARG:HH21	1.86	0.41
1:A:101:LEU:HB3	1:A:102:PRO:HD3	2.01	0.41
1:A:180:LEU:HB3	1:A:181:PRO:HD3	2.03	0.41
2:A:301:G2Q:HN3	2:A:301:G2Q:C3'	2.33	0.41
1:H:77:TRP:HA	1:H:158:MET:O	2.20	0.41
1:A:228:VAL:HG22	4:A:2171:HOH:O	2.21	0.41
1:C:44:ALA:HB3	1:C:58:TYR:CE1	2.56	0.41
1:F:80:VAL:O	1:F:81:TYR:C	2.63	0.41
1:G:142:MET:HE2	4:G:2040:HOH:O	2.20	0.41
1:G:174:PHE:CE2	2:G:301:G2Q:H2'	2.55	0.41
1:D:3:PRO:HG2	1:D:122:GLU:C	2.46	0.41
1:D:33:ASP:C	1:D:33:ASP:OD1	2.64	0.41
1:D:80:VAL:O	1:D:81:TYR:C	2.63	0.41
1:H:80:VAL:O	1:H:81:TYR:C	2.63	0.41
1:E:71:LYS:H	1:E:71:LYS:HG2	1.71	0.40
1:E:71:LYS:HG2	4:E:2082:HOH:O	2.20	0.40
1:E:73:LYS:HA	1:E:73:LYS:HD3	1.86	0.40
1:A:221:LYS:O	1:A:222:HIS:HB2	2.20	0.40
1:H:134:GLN:CD	1:H:134:GLN:H	2.29	0.40
1:D:198:LEU:HD22	1:D:247:GLU:HB2	2.04	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the sym-

metry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:154:GLU:OE1	1:F:31:ARG:NH1[1_565]	1.48	0.72
4:D:2183:HOH:O	4:F:2025:HOH:O[1_565]	2.12	0.08

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	250/259 (96%)	243 (97%)	7 (3%)	0	100	100
1	B	249/259 (96%)	247 (99%)	2 (1%)	0	100	100
1	C	250/259 (96%)	241 (96%)	9 (4%)	0	100	100
1	D	249/259 (96%)	242 (97%)	7 (3%)	0	100	100
1	E	249/259 (96%)	245 (98%)	4 (2%)	0	100	100
1	F	249/259 (96%)	241 (97%)	8 (3%)	0	100	100
1	G	250/259 (96%)	243 (97%)	7 (3%)	0	100	100
1	H	249/259 (96%)	240 (96%)	9 (4%)	0	100	100
All	All	1995/2072 (96%)	1942 (97%)	53 (3%)	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	220/226 (97%)	211 (96%)	9 (4%)	27	26
1	B	220/226 (97%)	213 (97%)	7 (3%)	34	35
1	C	220/226 (97%)	213 (97%)	7 (3%)	34	35
1	D	220/226 (97%)	211 (96%)	9 (4%)	27	26
1	E	220/226 (97%)	214 (97%)	6 (3%)	39	42
1	F	220/226 (97%)	205 (93%)	15 (7%)	14	11
1	G	220/226 (97%)	209 (95%)	11 (5%)	22	20
1	H	220/226 (97%)	212 (96%)	8 (4%)	31	31
All	All	1760/1808 (97%)	1688 (96%)	72 (4%)	27	26

All (72) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	30	VAL
1	A	35	SER
1	A	195	LEU
1	A	200	GLU
1	A	221	LYS
1	A	228	VAL
1	A	237	LEU
1	A	246	ARG
1	A	249	ARG
1	B	1	ILE
1	B	35	SER
1	B	39	LYS
1	B	129	LYS
1	B	195	LEU
1	B	232	ARG
1	B	237	LEU
1	C	1	ILE
1	C	30	VAL
1	C	92	ARG
1	C	123	LYS
1	C	195	LEU
1	C	200	GLU
1	C	221	LYS
1	D	30	VAL
1	D	92	ARG
1	D	115	ARG
1	D	117	ASN

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Mol	Chain	Res	Type
1	D	129	LYS
1	D	130	THR
1	D	159	VAL
1	D	216	LYS
1	D	228	VAL
1	E	1	ILE
1	E	30	VAL
1	E	39	LYS
1	E	195	LEU
1	E	237	LEU
1	E	246	ARG
1	F	1	ILE
1	F	20	ILE
1	F	30	VAL
1	F	33	ASP
1	F	73	LYS
1	F	92	ARG
1	F	115	ARG
1	F	123	LYS
1	F	127	ASP
1	F	129	LYS
1	F	186	LYS
1	F	195	LEU
1	F	221	LYS
1	F	232	ARG
1	F	246	ARG
1	G	30	VAL
1	G	31	ARG
1	G	92	ARG
1	G	108	SER
1	G	123	LYS
1	G	124	VAL
1	G	129	LYS
1	G	136	ARG
1	G	205	LYS
1	G	237	LEU
1	G	246	ARG
1	H	30	VAL
1	H	31	ARG
1	H	90	THR
1	H	123	LYS
1	H	124	VAL

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Mol	Chain	Res	Type
1	H	129	LYS
1	H	130	THR
1	H	246	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (25) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	66	GLN
1	A	89	ASN
1	A	117	ASN
1	A	242	ASN
1	B	49	ASN
1	B	89	ASN
1	B	242	ASN
1	C	66	GLN
1	C	89	ASN
1	C	117	ASN
1	C	242	ASN
1	D	89	ASN
1	D	117	ASN
1	D	242	ASN
1	E	89	ASN
1	E	117	ASN
1	E	242	ASN
1	F	89	ASN
1	G	66	GLN
1	G	89	ASN
1	G	242	ASN
1	H	89	ASN
1	H	117	ASN
1	H	134	GLN
1	H	242	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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

7 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
2	G2Q	C	301	-	39,39,39	3.65	14 (35%)	49,60,60	1.76	15 (30%)
2	G2Q	H	301	-	39,39,39	1.30	6 (15%)	49,60,60	1.71	6 (12%)
2	G2Q	B	301	-	39,39,39	2.78	16 (41%)	49,60,60	1.77	12 (24%)
3	NGD	F	5573	-	47,49,49	2.29	7 (14%)	66,75,75	2.22	18 (27%)
2	G2Q	E	301	-	39,39,39	2.60	15 (38%)	49,60,60	1.89	13 (26%)
2	G2Q	G	301	-	39,39,39	1.89	11 (28%)	49,60,60	2.26	14 (28%)
2	G2Q	A	301	-	39,39,39	2.82	17 (43%)	49,60,60	1.76	16 (32%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	G2Q	C	301	-	-	8/22/51/51	0/4/4/4
2	G2Q	H	301	-	-	9/22/51/51	0/4/4/4
2	G2Q	B	301	-	-	4/22/51/51	0/4/4/4
3	NGD	F	5573	-	-	7/30/62/62	0/5/5/5
2	G2Q	E	301	-	-	5/22/51/51	0/4/4/4
2	G2Q	G	301	-	-	6/22/51/51	0/4/4/4
2	G2Q	A	301	-	-	5/22/51/51	0/4/4/4

All (86) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	301	G2Q	PB-O3A	15.64	1.76	1.59
2	C	301	G2Q	PA-O3A	11.50	1.71	1.59
2	A	301	G2Q	PB-O3A	9.92	1.70	1.59
3	F	5573	NGD	C2-N1	8.27	1.44	1.35
2	A	301	G2Q	O4X-C1X	8.15	1.61	1.43
2	B	301	G2Q	O4X-C4X	7.88	1.57	1.44
2	E	301	G2Q	O4X-C4X	7.39	1.56	1.44
2	B	301	G2Q	PA-O3A	7.38	1.67	1.59
3	F	5573	NGD	O6A-C6A	7.07	1.37	1.23
2	B	301	G2Q	O4X-C1X	7.05	1.59	1.43
2	E	301	G2Q	O4X-C1X	6.23	1.57	1.43
2	C	301	G2Q	O4'-C1'	5.57	1.54	1.42
3	F	5573	NGD	O7-C7	5.52	1.34	1.24
2	A	301	G2Q	PA-O3A	5.21	1.65	1.59
2	C	301	G2Q	O4X-C4X	5.17	1.52	1.44
2	G	301	G2Q	PA-O3A	5.07	1.65	1.59
2	G	301	G2Q	PB-O3A	4.95	1.64	1.59
2	E	301	G2Q	PA-O5'	4.55	1.77	1.59
2	G	301	G2Q	C1'-N9	4.36	1.59	1.47
3	F	5573	NGD	PB-O3	4.34	1.64	1.59
2	E	301	G2Q	O4'-C1'	4.33	1.52	1.42
3	F	5573	NGD	PA-O3	4.27	1.64	1.59
2	A	301	G2Q	O4X-C4X	4.18	1.51	1.44
2	E	301	G2Q	C5'-C4'	4.06	1.63	1.51
3	F	5573	NGD	C6-N1	3.87	1.44	1.35
2	E	301	G2Q	C2X-C3X	3.77	1.60	1.53
2	B	301	G2Q	O4'-C1'	3.73	1.50	1.42
2	B	301	G2Q	PB-O3A	3.70	1.63	1.59
2	B	301	G2Q	PA-O5'	3.68	1.73	1.59
2	B	301	G2Q	C2X-C3X	3.64	1.59	1.53
2	E	301	G2Q	C2-N1	3.57	1.41	1.33
2	E	301	G2Q	C1X-C2X	3.52	1.57	1.51
2	B	301	G2Q	C5'-C4'	3.46	1.62	1.51
3	F	5573	NGD	O4'R-C1'R	3.42	1.45	1.40
2	H	301	G2Q	PB-O3A	3.40	1.63	1.59
2	E	301	G2Q	PB-O5X	3.38	1.72	1.59
2	G	301	G2Q	O4'-C1'	3.30	1.49	1.42
2	A	301	G2Q	PB-O1B	3.23	1.62	1.50
2	B	301	G2Q	PB-O5X	3.19	1.71	1.59
2	A	301	G2Q	O4'-C1'	3.18	1.49	1.42
2	C	301	G2Q	C1'-N9	3.07	1.56	1.47
2	G	301	G2Q	C2-N1	3.02	1.40	1.33
2	E	301	G2Q	C6-N1	2.98	1.43	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	301	G2Q	C1X-C2X	2.86	1.56	1.51
2	B	301	G2Q	PB-O1B	2.84	1.60	1.50
2	C	301	G2Q	C2-N1	2.83	1.40	1.33
2	A	301	G2Q	C2X-C3X	2.70	1.58	1.53
2	B	301	G2Q	C6-N1	2.67	1.42	1.38
2	A	301	G2Q	PB-O5X	2.64	1.69	1.59
2	E	301	G2Q	O2'-C2'	2.56	1.49	1.43
2	C	301	G2Q	C5'-C4'	2.54	1.59	1.51
2	C	301	G2Q	PB-O5X	2.51	1.69	1.59
2	A	301	G2Q	C2-N1	2.51	1.39	1.33
2	C	301	G2Q	C6-N1	2.46	1.42	1.38
2	B	301	G2Q	C2-N1	2.45	1.39	1.33
2	A	301	G2Q	O4'-C4'	2.44	1.50	1.45
2	A	301	G2Q	O5X-C5X	2.38	1.53	1.44
2	H	301	G2Q	C5-N7	-2.37	1.34	1.39
2	A	301	G2Q	C6-N1	2.37	1.42	1.38
2	A	301	G2Q	C5'-C4'	2.36	1.58	1.51
2	H	301	G2Q	PA-O3A	2.35	1.62	1.59
2	B	301	G2Q	C5-N7	-2.32	1.34	1.39
2	B	301	G2Q	C5-C4	2.32	1.42	1.38
2	G	301	G2Q	C4-N9	-2.29	1.34	1.37
2	H	301	G2Q	PB-O5X	2.29	1.68	1.59
2	B	301	G2Q	C1'-N9	2.29	1.54	1.47
2	C	301	G2Q	PA-O5'	2.28	1.68	1.59
2	E	301	G2Q	C5-C4	2.25	1.42	1.38
2	A	301	G2Q	PA-O2A	-2.24	1.45	1.55
2	A	301	G2Q	PA-O5'	2.22	1.68	1.59
2	E	301	G2Q	PA-O3A	2.21	1.61	1.59
2	E	301	G2Q	O5X-C5X	2.21	1.53	1.44
2	A	301	G2Q	C5-C4	2.20	1.42	1.38
2	C	301	G2Q	PA-O2A	-2.10	1.45	1.55
2	C	301	G2Q	C3'-C4'	2.10	1.58	1.53
2	H	301	G2Q	C5X-C4X	2.09	1.57	1.51
2	H	301	G2Q	C6-C5	-2.08	1.39	1.45
2	G	301	G2Q	O4'-C4'	2.08	1.49	1.45
2	G	301	G2Q	C2-N3	-2.07	1.32	1.37
2	G	301	G2Q	PB-O5X	2.05	1.67	1.59
2	E	301	G2Q	C4-N9	-2.04	1.35	1.37
2	B	301	G2Q	C2-N3	-2.04	1.32	1.37
2	G	301	G2Q	C6-N1	2.03	1.41	1.38
2	C	301	G2Q	O4X-C1X	2.01	1.48	1.43
2	A	301	G2Q	O6-C6	2.01	1.27	1.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	G	301	G2Q	PA-O2A	-2.00	1.46	1.55

All (94) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	F	5573	NGD	C6-N1-C2	-9.51	113.79	121.88
2	G	301	G2Q	C5-C4-N9	9.12	110.47	106.25
3	F	5573	NGD	C5A-C4A-N3A	-6.34	118.30	128.39
2	G	301	G2Q	C2'-C1'-N9	6.21	130.53	113.25
3	F	5573	NGD	C2A-N3A-C4A	5.95	122.55	112.30
2	H	301	G2Q	N2-C2-N3	5.57	128.51	116.76
2	H	301	G2Q	C5-C4-N9	-5.49	103.70	106.25
2	G	301	G2Q	C5'-C4'-C3'	-4.98	97.28	115.21
2	E	301	G2Q	C1X-C2X-C3X	4.85	109.39	101.63
2	C	301	G2Q	O4'-C1'-N9	4.83	119.30	108.36
2	B	301	G2Q	C5X-C4X-C3X	-4.75	98.11	115.21
3	F	5573	NGD	N9A-C4A-N3A	4.52	135.00	125.95
2	B	301	G2Q	O3A-PA-O1A	-4.29	97.79	110.70
2	H	301	G2Q	N2-C2-N1	-3.93	111.99	119.67
2	A	301	G2Q	O3A-PA-O1A	-3.74	99.45	110.70
2	E	301	G2Q	O3A-PA-O1A	-3.62	99.83	110.70
2	E	301	G2Q	O2B-PB-O1B	3.56	129.00	112.44
2	G	301	G2Q	C4-C5-N7	-3.55	104.65	109.26
2	E	301	G2Q	O4X-C1X-C2X	-3.55	99.63	105.76
2	B	301	G2Q	C1X-C2X-C3X	3.52	107.25	101.63
2	C	301	G2Q	O4'-C4'-C5'	-3.44	98.32	109.33
3	F	5573	NGD	C3-C7-N7	3.40	121.93	117.74
2	E	301	G2Q	C5X-C4X-C3X	-3.37	103.07	115.21
2	H	301	G2Q	C8-N9-C4	3.32	109.47	106.67
2	E	301	G2Q	O3X-C3X-C2X	3.24	119.64	111.97
2	A	301	G2Q	N2-C2-N3	3.15	123.41	116.76
2	B	301	G2Q	C8-N7-C5	3.12	109.81	104.26
2	B	301	G2Q	O2B-PB-O1B	3.05	126.65	112.44
3	F	5573	NGD	C5-C6-N1	3.03	124.52	120.38
2	A	301	G2Q	C8-N7-C5	3.00	109.61	104.26
2	G	301	G2Q	O4X-C4X-C3X	2.98	107.39	104.63
2	E	301	G2Q	O4'-C4'-C3'	-2.97	99.26	105.15
2	A	301	G2Q	C4-C5-N7	-2.96	105.42	109.26
2	C	301	G2Q	O4X-C4X-C3X	2.93	107.34	104.63
2	C	301	G2Q	C8-N9-C4	2.89	109.11	106.67
2	A	301	G2Q	O6-C6-N1	-2.88	115.43	120.27
2	A	301	G2Q	O2A-PA-O3A	-2.86	99.53	107.27

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	301	G2Q	O4X-C1X-C2X	-2.86	100.83	105.76
2	E	301	G2Q	C8-N7-C5	2.85	109.35	104.26
2	C	301	G2Q	C5'-C4'-C3'	2.85	125.48	115.21
2	E	301	G2Q	O2B-PB-O3A	-2.85	99.57	107.27
3	F	5573	NGD	C3'-C2'-C1'	2.82	106.80	101.46
2	A	301	G2Q	N3-C2-N1	-2.80	118.20	123.32
2	G	301	G2Q	C8-N7-C5	2.76	109.18	104.26
2	C	301	G2Q	C1X-C2X-C3X	2.76	106.04	101.63
2	G	301	G2Q	C1'-N9-C8	2.75	134.53	126.73
3	F	5573	NGD	O4'-C1'-N9A	2.74	114.57	108.36
2	A	301	G2Q	O4X-C1X-C2X	-2.73	101.06	105.76
2	B	301	G2Q	O2A-PA-O3A	-2.66	100.09	107.27
2	C	301	G2Q	C8-N7-C5	2.65	108.98	104.26
2	A	301	G2Q	O2B-PB-O1B	2.65	124.75	112.44
2	C	301	G2Q	C2'-C1'-N9	-2.60	106.01	113.25
3	F	5573	NGD	O6A-C6A-C5A	-2.53	119.86	126.53
3	F	5573	NGD	C6-N1-C1'R	-2.53	114.77	119.73
2	A	301	G2Q	C1X-C2X-C3X	2.51	105.65	101.63
2	C	301	G2Q	O2A-PA-O1A	2.46	123.90	112.44
2	G	301	G2Q	O6-C6-N1	-2.45	116.15	120.27
2	G	301	G2Q	C3'-C2'-C1'	2.41	106.02	101.46
2	A	301	G2Q	O3'-C3'-C2'	-2.38	104.20	111.82
2	E	301	G2Q	N9-C8-N7	-2.37	109.00	113.40
3	F	5573	NGD	C5A-C6A-N1A	2.36	119.26	113.25
3	F	5573	NGD	C2A-N1A-C6A	-2.34	120.87	125.11
2	B	301	G2Q	O2A-PA-O1A	2.34	123.31	112.44
2	G	301	G2Q	O4'-C1'-C2'	-2.33	101.63	106.62
2	B	301	G2Q	N9-C8-N7	-2.33	109.08	113.40
2	H	301	G2Q	O4'-C1'-C2'	-2.32	101.65	106.62
2	G	301	G2Q	C6-C5-N7	2.32	137.67	130.12
3	F	5573	NGD	C2'-C1'-N9A	-2.32	106.80	113.25
3	F	5573	NGD	C4'R-O4'R-C1'R	2.31	112.04	109.92
2	C	301	G2Q	N9-C8-N7	-2.29	109.15	113.40
2	A	301	G2Q	N9-C8-N7	-2.28	109.18	113.40
2	C	301	G2Q	O3X-C3X-C4X	-2.27	104.55	111.08
2	B	301	G2Q	N3-C2-N1	-2.27	119.16	123.32
2	A	301	G2Q	C8-N9-C4	2.27	108.59	106.67
2	G	301	G2Q	O2A-PA-O5'	-2.22	97.52	107.57
2	E	301	G2Q	C4-C5-N7	-2.21	106.39	109.26
2	C	301	G2Q	O2B-PB-O1B	2.20	122.70	112.44
2	A	301	G2Q	O2X-C2X-C3X	2.19	115.74	111.43
2	C	301	G2Q	N2-C2-N3	2.15	121.30	116.76

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	301	G2Q	C8-N9-C4	2.15	108.48	106.67
2	A	301	G2Q	C6-N1-C2	2.13	125.50	119.24
2	C	301	G2Q	O3'-C3'-C4'	2.09	117.09	111.08
2	G	301	G2Q	N9-C8-N7	-2.09	109.52	113.40
2	H	301	G2Q	N3-C2-N1	-2.09	119.50	123.32
2	G	301	G2Q	O4'-C4'-C5'	2.07	115.97	109.33
3	F	5573	NGD	O7-C7-N7	-2.07	119.62	122.62
3	F	5573	NGD	C2'R-C3'R-C4'R	2.06	106.59	102.61
2	C	301	G2Q	O3'-C3'-C2'	-2.05	105.24	111.82
2	A	301	G2Q	C5-C4-N9	2.04	107.19	106.25
2	B	301	G2Q	C4-C5-N7	-2.04	106.62	109.26
2	B	301	G2Q	O2X-C2X-C3X	-2.03	107.42	111.43
3	F	5573	NGD	O1A-PA-O2A	2.03	121.89	112.44
2	E	301	G2Q	N2-C2-N3	2.03	121.04	116.76
3	F	5573	NGD	N9A-C8A-N7A	-2.02	109.66	113.40

There are no chirality outliers.

All (44) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	301	G2Q	C5'-O5'-PA-O1A
2	C	301	G2Q	C5'-O5'-PA-O3A
2	G	301	G2Q	C5'-O5'-PA-O1A
2	H	301	G2Q	C5'-O5'-PA-O1A
2	H	301	G2Q	C5'-O5'-PA-O2A
2	H	301	G2Q	C5'-O5'-PA-O3A
3	F	5573	NGD	O4'-C1'-N9A-C8A
3	F	5573	NGD	O4'-C1'-N9A-C4A
3	F	5573	NGD	O4'-C4'-C5'-O5'
3	F	5573	NGD	C2'R-C1'R-N1-C6
2	H	301	G2Q	O4'-C4'-C5'-O5'
3	F	5573	NGD	C3'-C4'-C5'-O5'
2	E	301	G2Q	C3'-C4'-C5'-O5'
2	H	301	G2Q	C3'-C4'-C5'-O5'
2	G	301	G2Q	C3'-C4'-C5'-O5'
2	B	301	G2Q	C3'-C4'-C5'-O5'
2	G	301	G2Q	O4'-C4'-C5'-O5'
2	E	301	G2Q	O4'-C4'-C5'-O5'
2	H	301	G2Q	PA-O3A-PB-O1B
2	C	301	G2Q	PB-O3A-PA-O5'
2	E	301	G2Q	PB-O3A-PA-O5'
2	H	301	G2Q	PB-O3A-PA-O5'

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Mol	Chain	Res	Type	Atoms
3	F	5573	NGD	PB-O3-PA-O5'
2	B	301	G2Q	O4'-C4'-C5'-O5'
2	A	301	G2Q	C3'-C4'-C5'-O5'
2	B	301	G2Q	C4X-C5X-O5X-PB
2	G	301	G2Q	C4X-C5X-O5X-PB
2	C	301	G2Q	C5'-O5'-PA-O2A
2	C	301	G2Q	C4X-C5X-O5X-PB
2	E	301	G2Q	C4X-C5X-O5X-PB
2	G	301	G2Q	C4'-C5'-O5'-PA
2	H	301	G2Q	C4X-C5X-O5X-PB
2	H	301	G2Q	PA-O3A-PB-O2B
2	A	301	G2Q	O4'-C4'-C5'-O5'
3	F	5573	NGD	C4'R-C5'R-O5'R-PB
2	B	301	G2Q	PB-O3A-PA-O5'
2	A	301	G2Q	C4X-C5X-O5X-PB
2	A	301	G2Q	C4'-C5'-O5'-PA
2	C	301	G2Q	C4'-C5'-O5'-PA
2	C	301	G2Q	C2'-C1'-N9-C4
2	G	301	G2Q	PA-O3A-PB-O2B
2	E	301	G2Q	C2'-C1'-N9-C8
2	A	301	G2Q	PA-O3A-PB-O2B
2	C	301	G2Q	O4'-C4'-C5'-O5'

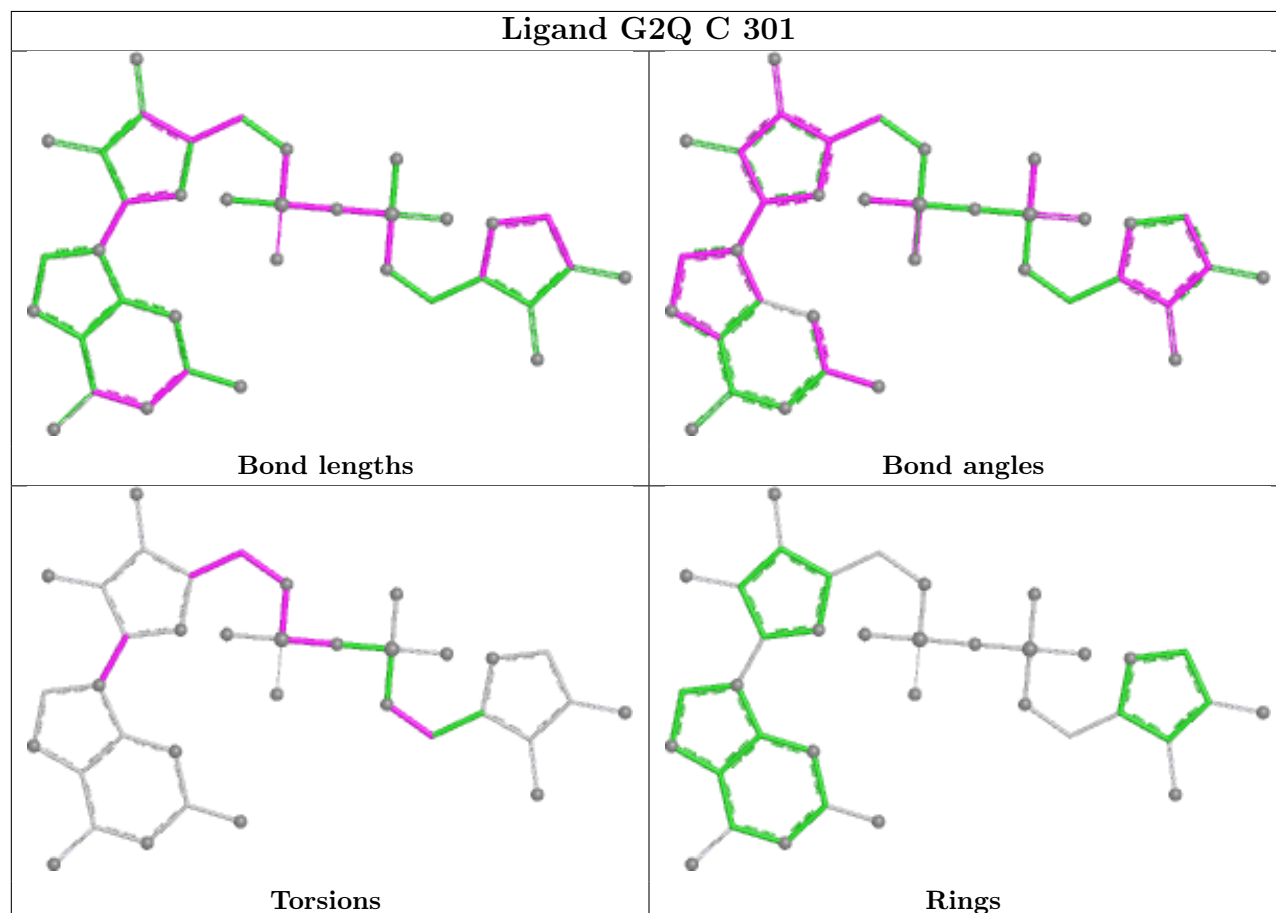
There are no ring outliers.

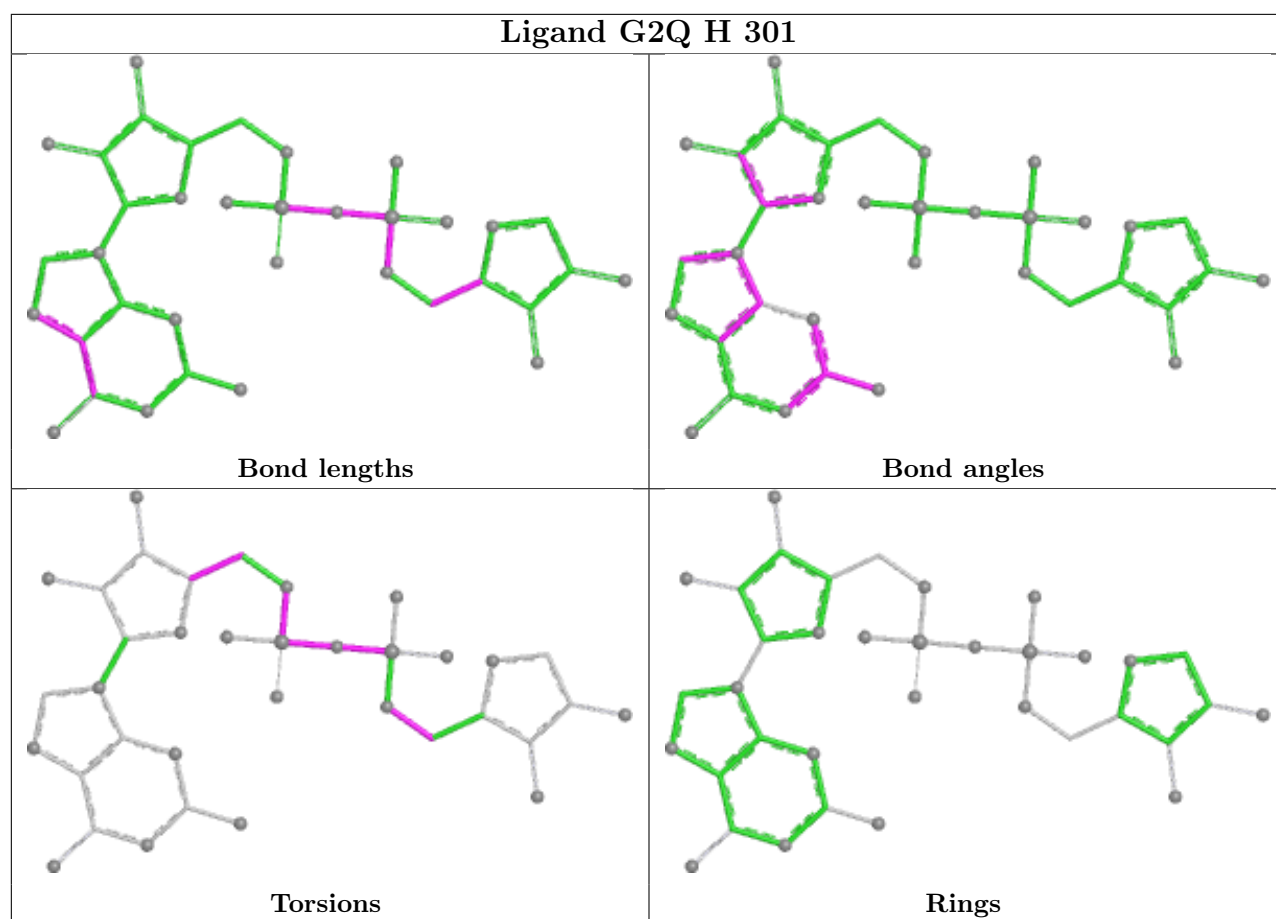
7 monomers are involved in 65 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	301	G2Q	13	0
2	H	301	G2Q	5	0
2	B	301	G2Q	12	0
3	F	5573	NGD	5	0
2	E	301	G2Q	11	0
2	G	301	G2Q	9	0
2	A	301	G2Q	10	0

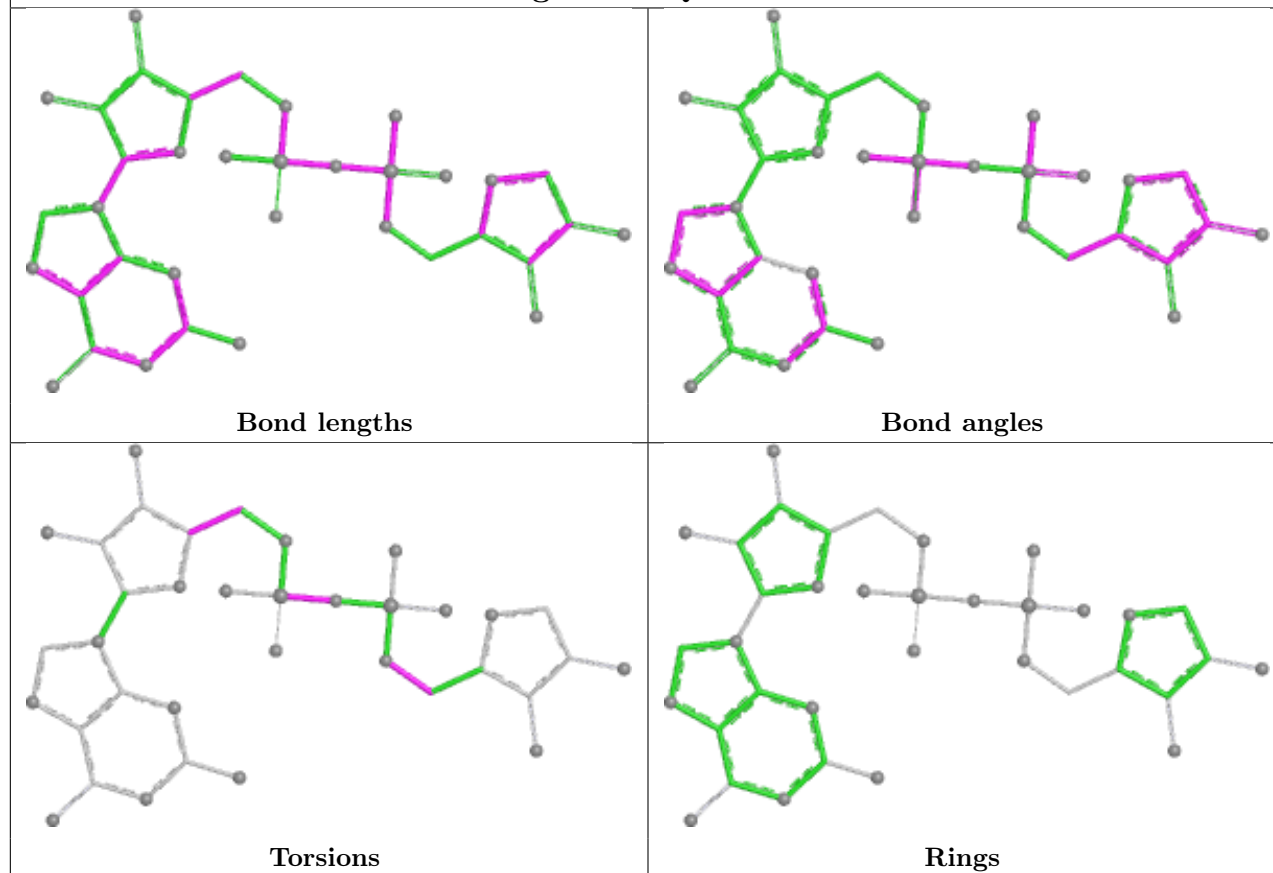
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

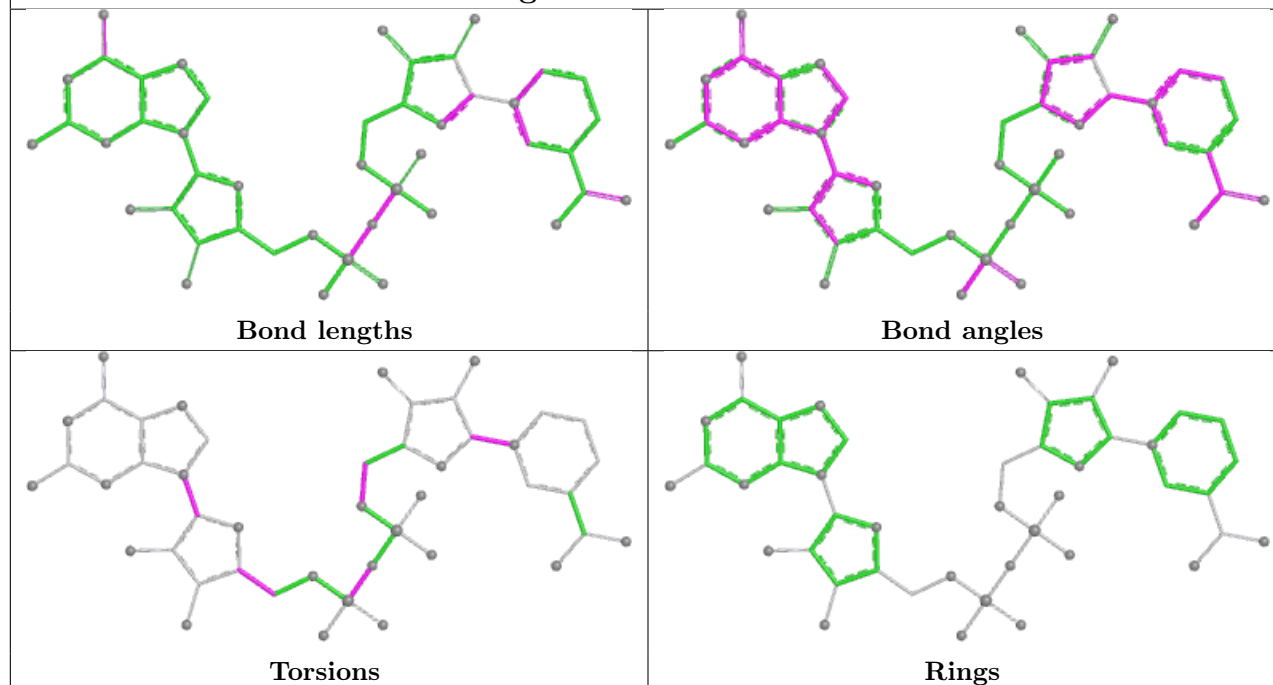


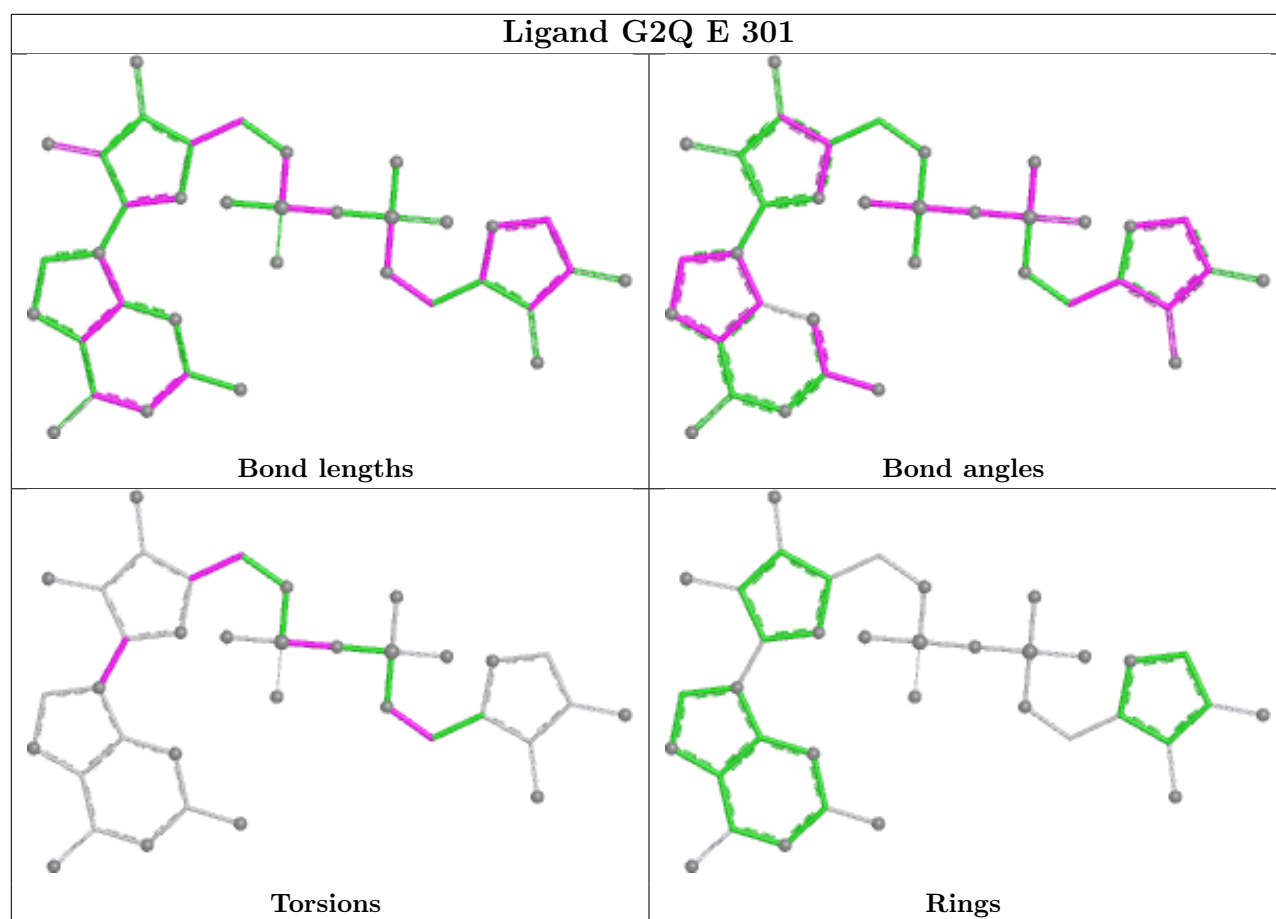


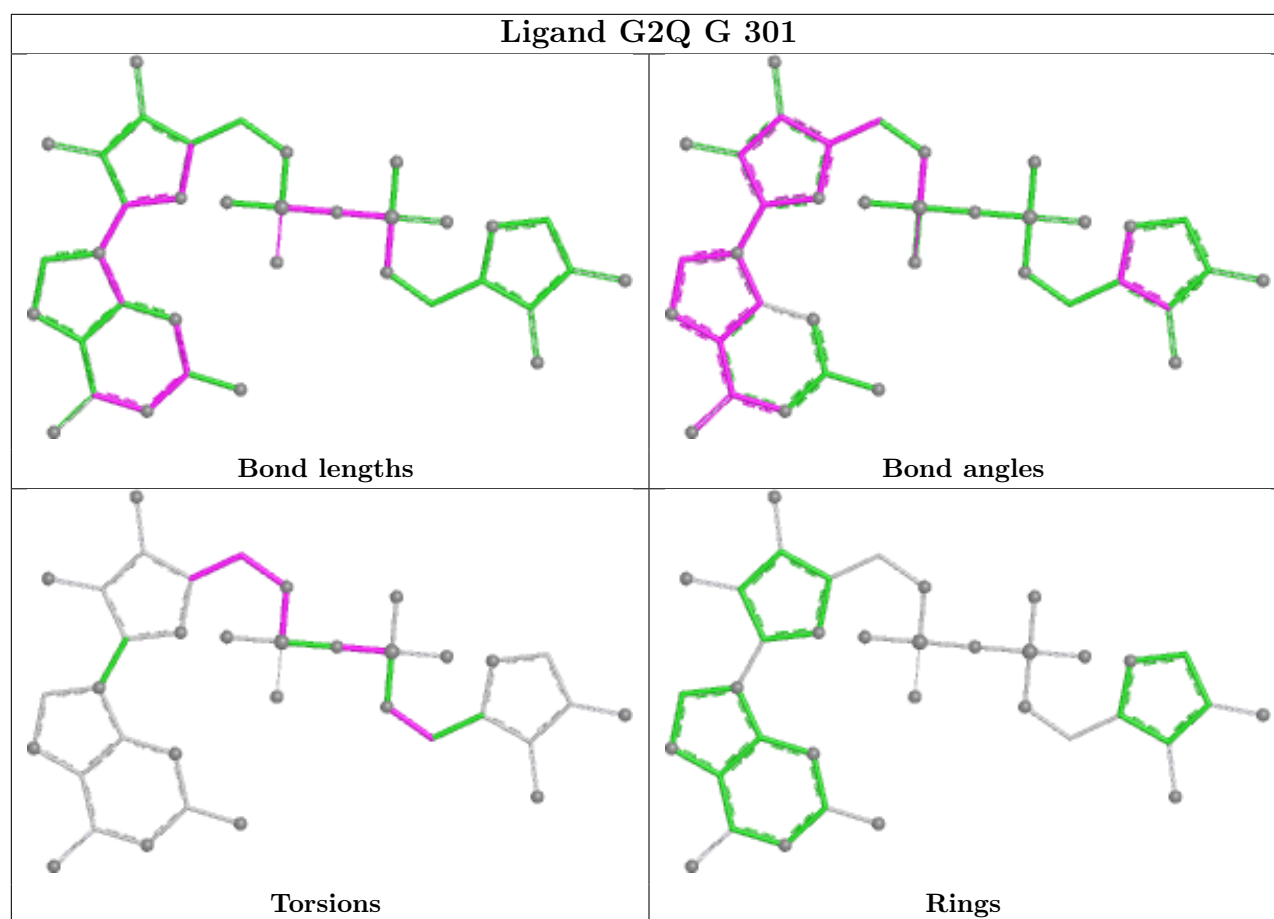
Ligand G2Q B 301

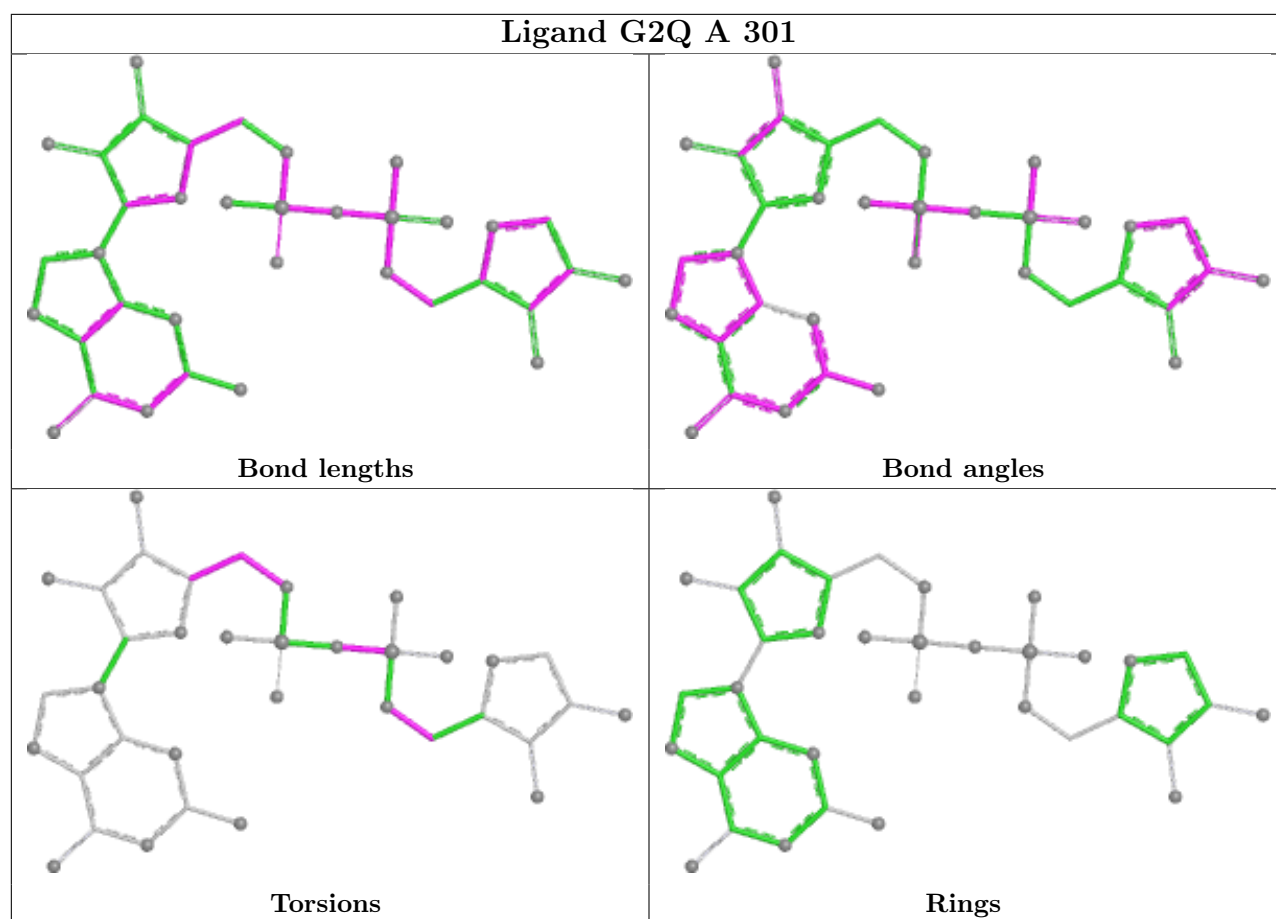


Ligand NGD F 5573









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	252/259 (97%)	-0.32	3 (1%) 76 76	15, 30, 48, 57	0
1	B	251/259 (96%)	-0.32	5 (1%) 65 64	19, 31, 49, 61	0
1	C	252/259 (97%)	-0.31	3 (1%) 76 76	17, 29, 48, 57	0
1	D	251/259 (96%)	-0.14	3 (1%) 76 76	18, 32, 53, 76	0
1	E	251/259 (96%)	-0.23	5 (1%) 65 64	19, 31, 47, 60	0
1	F	251/259 (96%)	-0.12	3 (1%) 76 76	19, 36, 56, 75	0
1	G	252/259 (97%)	0.12	4 (1%) 70 70	29, 46, 62, 76	0
1	H	251/259 (96%)	0.08	2 (0%) 82 82	30, 45, 63, 74	0
All	All	2011/2072 (97%)	-0.16	28 (1%) 73 73	15, 35, 56, 76	0

All (28) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	128	PHE	4.9
1	H	128	PHE	3.9
1	C	0	ALA	3.8
1	H	174	PHE	3.8
1	C	174	PHE	3.8
1	G	174	PHE	3.7
1	A	174	PHE	3.7
1	E	174	PHE	3.6
1	D	128	PHE	3.4
1	G	128	PHE	3.4
1	F	174	PHE	3.2
1	F	128	PHE	3.0
1	A	0	ALA	2.9
1	B	174	PHE	2.9
1	B	128	PHE	2.7
1	D	174	PHE	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	128	PHE	2.5
1	B	52	ASP	2.3
1	E	251	ALA	2.3
1	G	251	ALA	2.3
1	B	251	ALA	2.2
1	D	251	ALA	2.2
1	G	127	ASP	2.2
1	E	249	ARG	2.1
1	E	246	ARG	2.1
1	B	246	ARG	2.1
1	C	246	ARG	2.1
1	F	1	ILE	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](#)

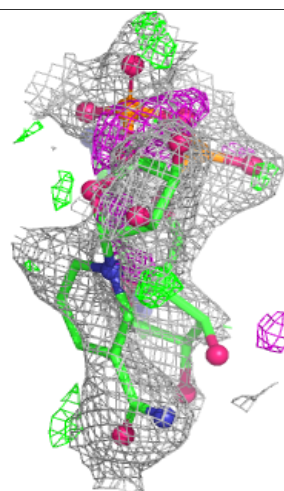
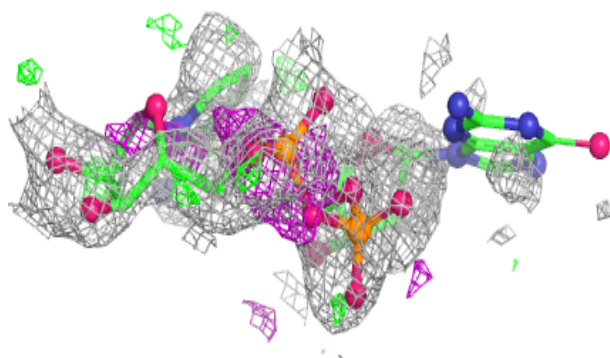
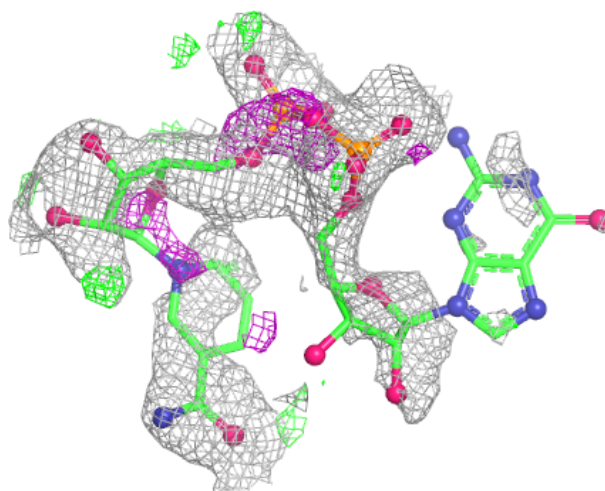
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	NGD	F	5573	45/45	0.72	0.16	64,99,123,125	0
2	G2Q	E	301	36/36	0.82	0.14	27,51,57,59	0
2	G2Q	A	301	36/36	0.82	0.13	20,43,54,56	0
2	G2Q	B	301	36/36	0.83	0.12	31,48,54,56	0
2	G2Q	C	301	36/36	0.83	0.13	27,50,61,62	0
2	G2Q	H	301	36/36	0.86	0.14	58,72,78,81	0
2	G2Q	G	301	36/36	0.86	0.12	60,69,78,81	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.

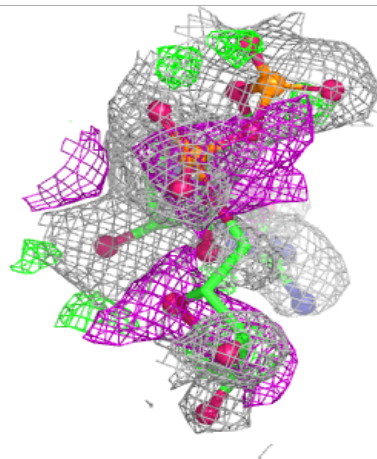
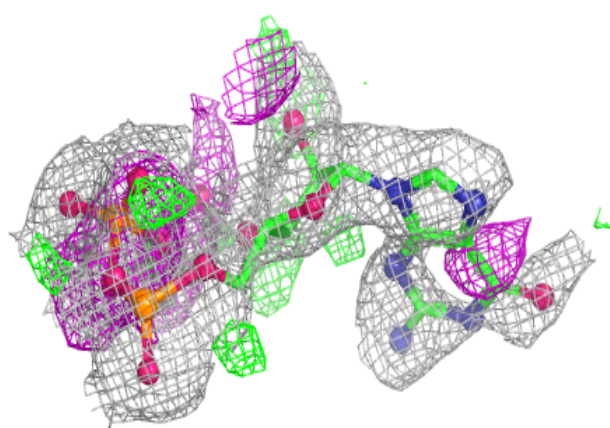
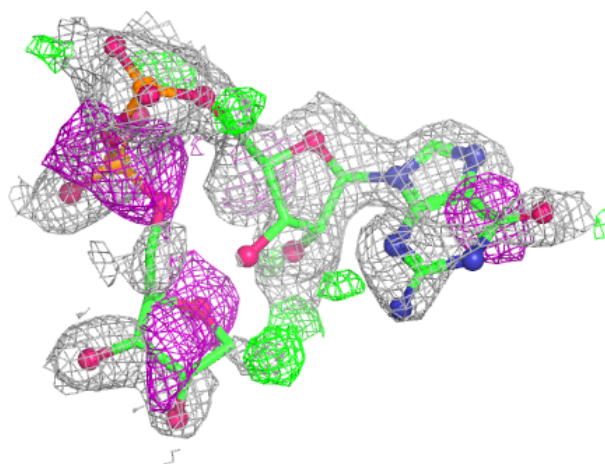
Electron density around NGD F 5573:

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



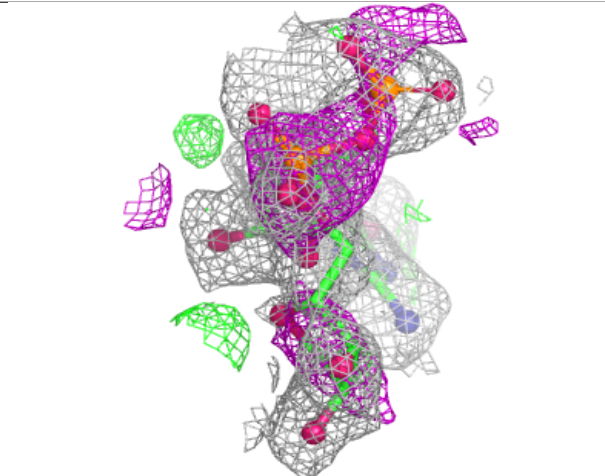
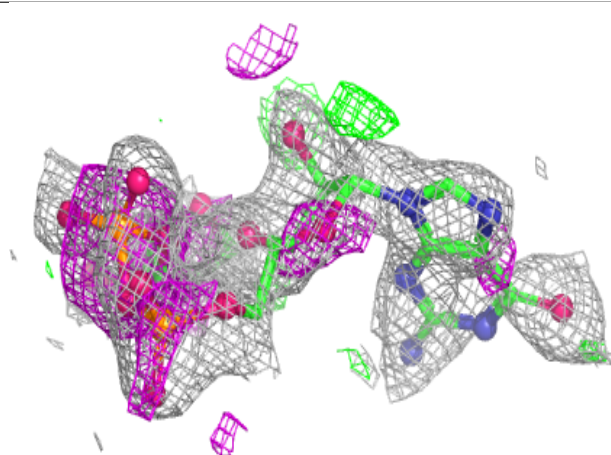
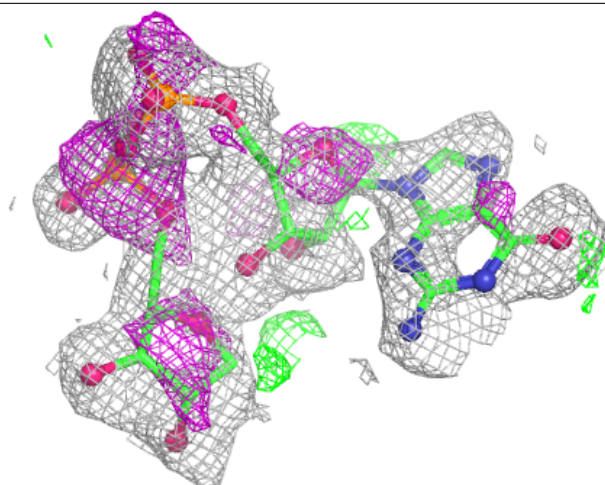
Electron density around G2Q E 301:

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



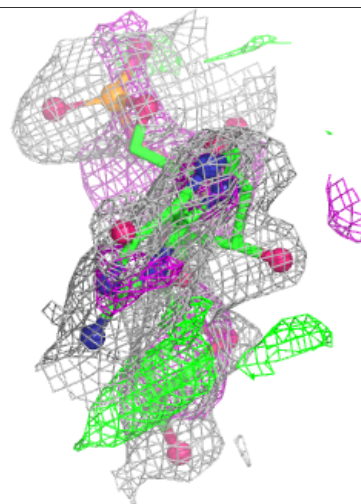
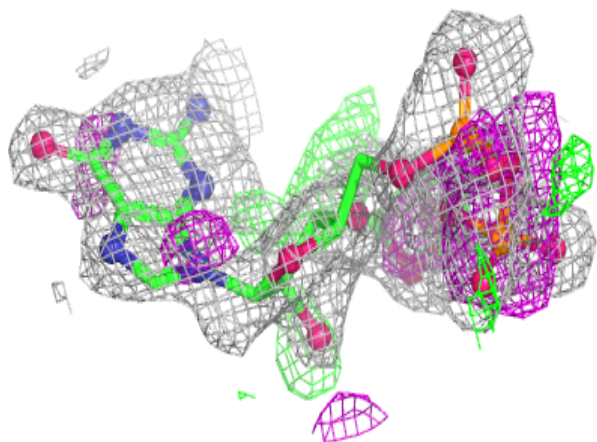
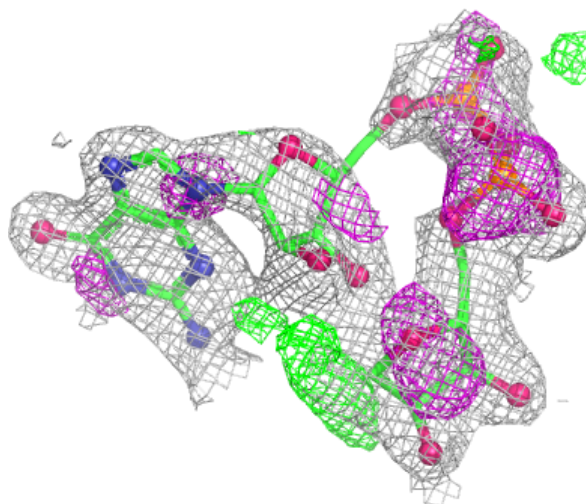
Electron density around G2Q A 301:

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



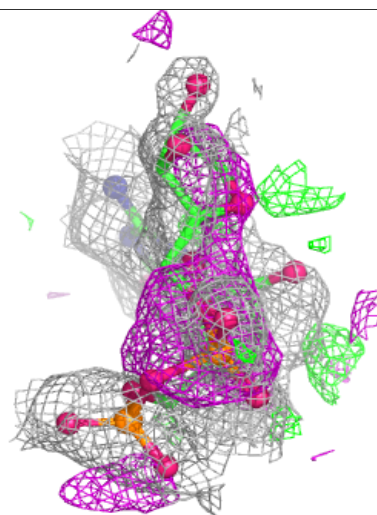
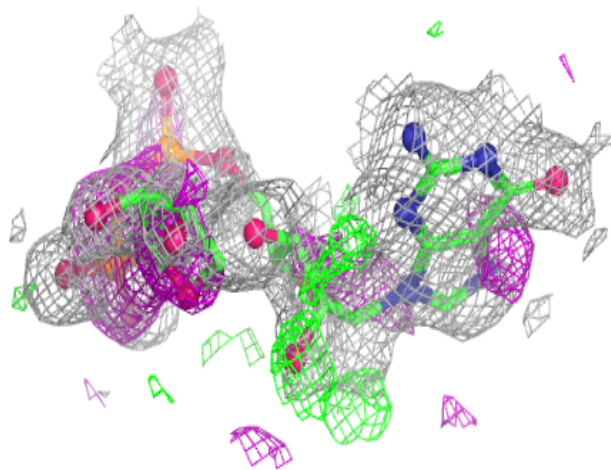
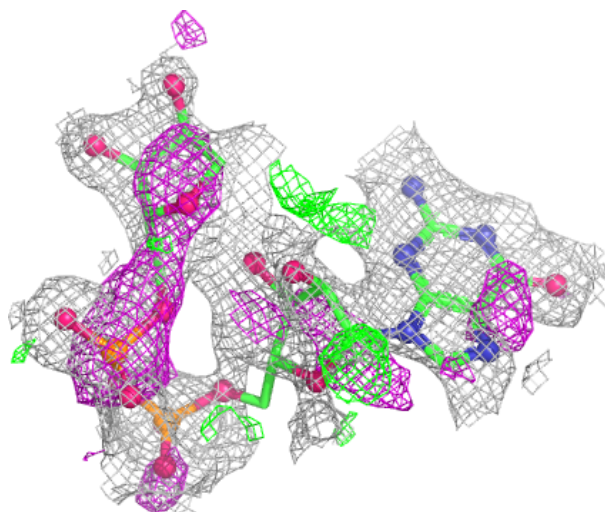
Electron density around G2Q B 301:

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



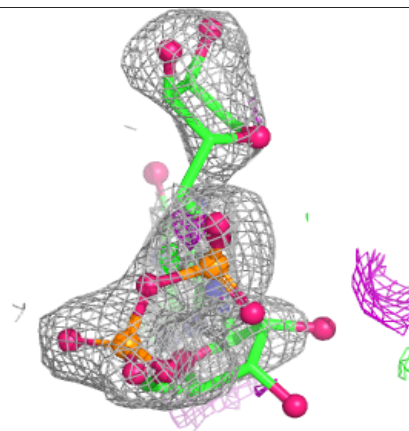
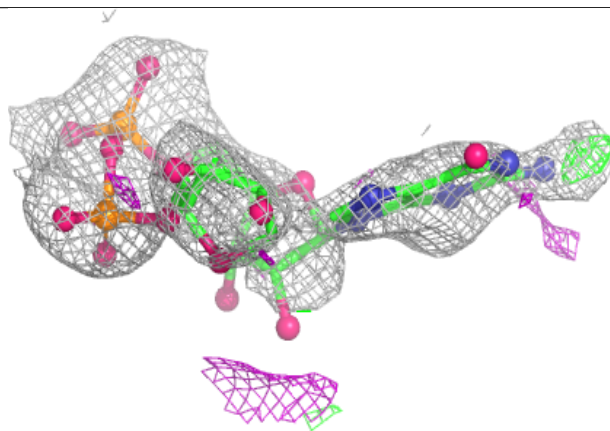
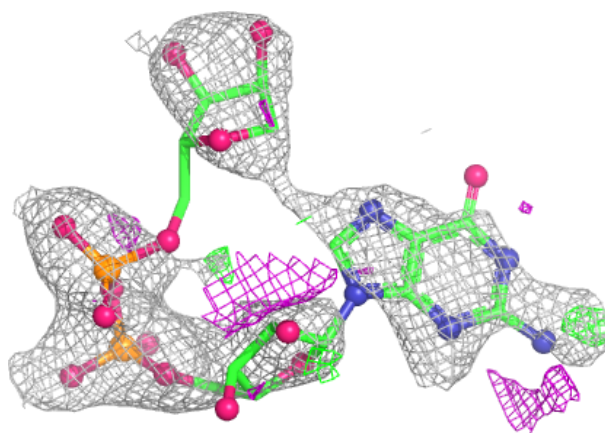
Electron density around G2Q C 301:

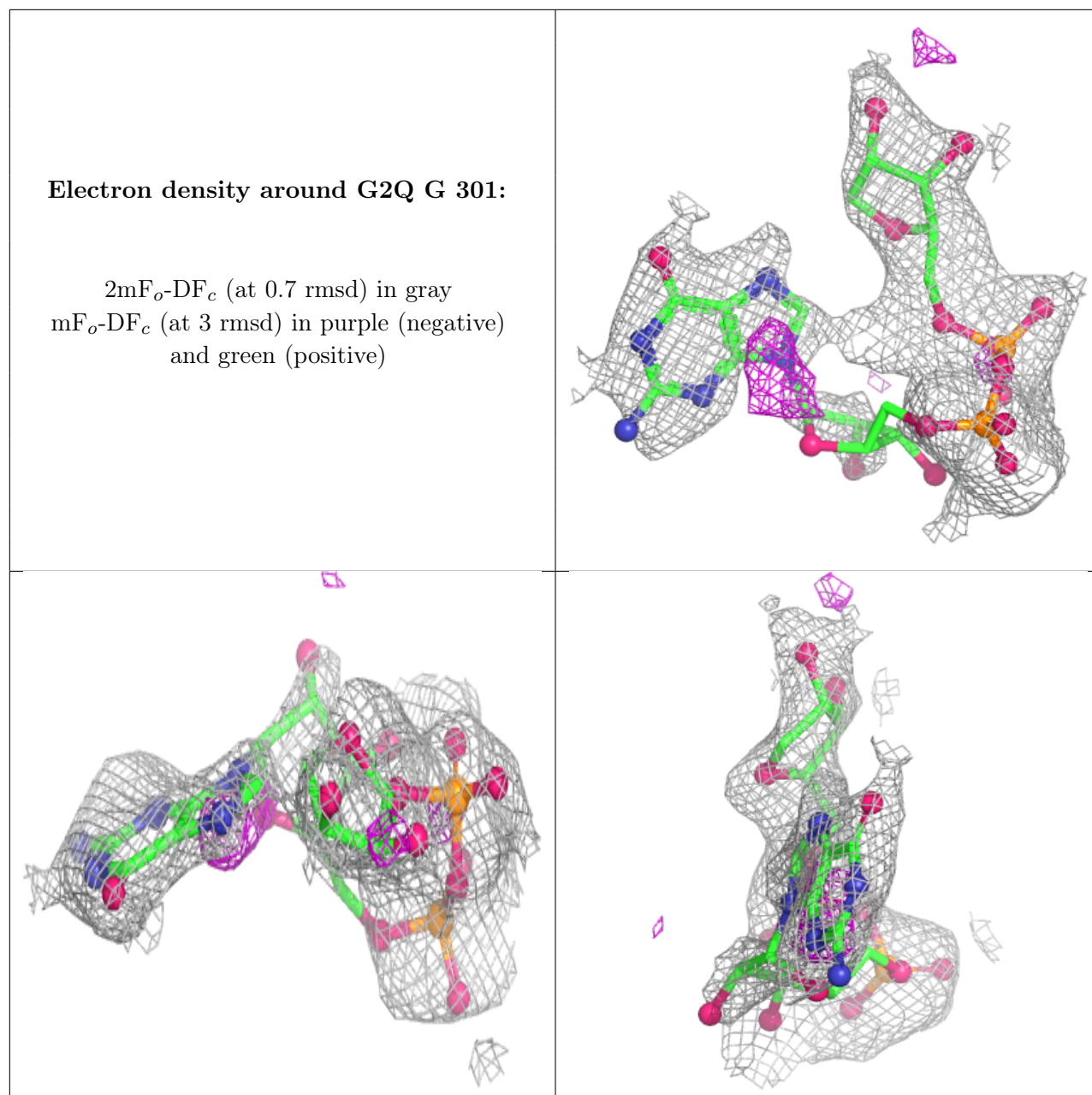
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around G2Q H 301:

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





6.5 Other polymers ⓘ

There are no such residues in this entry.