



Full wwPDB X-ray Structure Validation Report ⓘ

May 26, 2020 – 08:39 am BST

PDB ID : 4LUC
Title : Crystal Structure of small molecule disulfide 6 bound to K-Ras G12C
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Deposited on : 2013-07-25
Resolution : 1.29 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

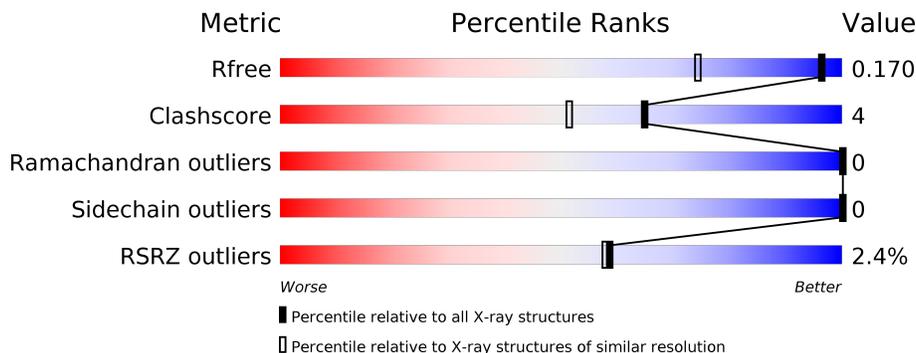
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.29 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	1058 (1.30-1.30)
Clashscore	141614	1101 (1.30-1.30)
Ramachandran outliers	138981	1058 (1.30-1.30)
Sidechain outliers	138945	1058 (1.30-1.30)
RSRZ outliers	127900	1029 (1.30-1.30)

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	170	
1	B	170	

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 5925 atoms, of which 2773 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 GTPase KRas.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace	
			Total	C	H	N	O				S
1	A	168	2715	855	1357	231	267	5	0	5	0
1	B	168	2705	852	1350	231	266	6	0	4	0

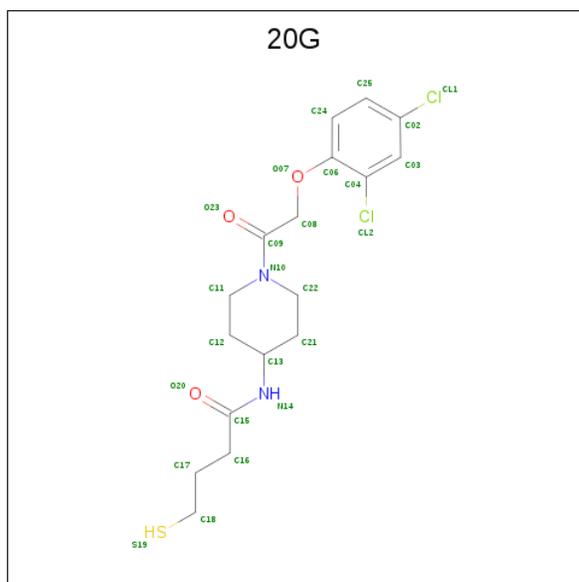
There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	GLY	-	EXPRESSION TAG	UNP P01116
A	12	CYS	GLY	VARIANT	UNP P01116
A	51	SER	CYS	ENGINEERED MUTATION	UNP P01116
A	80	LEU	CYS	ENGINEERED MUTATION	UNP P01116
A	118	SER	CYS	ENGINEERED MUTATION	UNP P01116
A	151	GLY	ARG	SEE REMARK 999	UNP P01116
A	153	ASP	GLU	SEE REMARK 999	UNP P01116
A	165	LYS	GLN	SEE REMARK 999	UNP P01116
A	166	HIS	TYR	SEE REMARK 999	UNP P01116
A	167	LYS	ARG	SEE REMARK 999	UNP P01116
A	168	GLU	LEU	SEE REMARK 999	UNP P01116
B	0	GLY	-	EXPRESSION TAG	UNP P01116
B	12	CYS	GLY	VARIANT	UNP P01116
B	51	SER	CYS	ENGINEERED MUTATION	UNP P01116
B	80	LEU	CYS	ENGINEERED MUTATION	UNP P01116
B	118	SER	CYS	ENGINEERED MUTATION	UNP P01116
B	151	GLY	ARG	SEE REMARK 999	UNP P01116
B	153	ASP	GLU	SEE REMARK 999	UNP P01116
B	165	LYS	GLN	SEE REMARK 999	UNP P01116
B	166	HIS	TYR	SEE REMARK 999	UNP P01116
B	167	LYS	ARG	SEE REMARK 999	UNP P01116
B	168	GLU	LEU	SEE REMARK 999	UNP P01116

- Molecule 2 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	2	Total Ca 2 2	0	0
2	A	2	Total Ca 2 2	0	0

- Molecule 3 is N-{1-[(2,4-dichlorophenoxy)acetyl]piperidin-4-yl}-4-sulfanylbutanamide (three-letter code: 20G) (formula: $C_{17}H_{22}Cl_2N_2O_3S$).



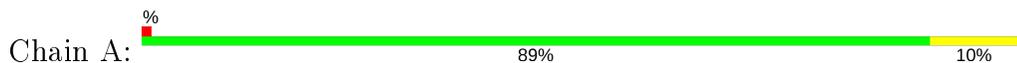
Mol	Chain	Residues	Atoms							ZeroOcc	AltConf
			Total	C	Cl	H	N	O	S		
3	A	1	Total 46	C 17	Cl 2	H 21	N 2	O 3	S 1	0	0
3	B	1	Total 46	C 17	Cl 2	H 21	N 2	O 3	S 1	0	0

- Molecule 4 is GUANOSINE-5'-DIPHOSPHATE (three-letter code: GDP) (formula: $C_{10}H_{15}N_5O_{11}P_2$).

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: GTPase KRas



- Molecule 1: GTPase KRas



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	32.82Å 39.08Å 62.46Å 77.86° 81.54° 77.55°	Depositor
Resolution (Å)	19.37 – 1.29 19.37 – 1.29	Depositor EDS
% Data completeness (in resolution range)	95.0 (19.37-1.29) 92.0 (19.37-1.29)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.60 (at 1.29Å)	Xtrriage
Refinement program	PHENIX dev_1402	Depositor
R, R_{free}	0.152 , 0.169 0.152 , 0.170	Depositor DCC
R_{free} test set	2000 reflections (2.83%)	wwPDB-VP
Wilson B-factor (Å ²)	11.7	Xtrriage
Anisotropy	0.365	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.47 , 49.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.98	EDS
Total number of atoms	5925	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.28% 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: GDP, CA, 20G

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	1.03	0/1394	1.01	5/1882 (0.3%)
1	B	0.90	1/1388 (0.1%)	0.97	5/1873 (0.3%)
All	All	0.97	1/2782 (0.0%)	0.99	10/3755 (0.3%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	157	TYR	CE1-CZ	-5.36	1.31	1.38

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	102	ARG	NE-CZ-NH2	-11.24	114.68	120.30
1	B	33	ASP	CB-CG-OD1	7.26	124.84	118.30
1	A	102	ARG	NE-CZ-NH2	-7.03	116.78	120.30
1	A	79	LEU	CB-CG-CD2	5.83	120.91	111.00
1	B	111[A]	MET	CG-SD-CE	-5.61	91.22	100.20
1	B	111[B]	MET	CG-SD-CE	-5.61	91.22	100.20
1	B	97	ARG	NE-CZ-NH1	5.58	123.09	120.30
1	A	99	GLN	CA-CB-CG	5.54	125.59	113.40
1	A	137	TYR	CB-CG-CD2	-5.47	117.72	121.00
1	A	92	ASP	CB-CG-OD1	5.05	122.84	118.30

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	1358	1357	1355	18	0
1	B	1355	1350	1348	5	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
3	A	25	21	21	0	0
3	B	25	21	21	0	0
4	A	28	12	12	0	0
4	B	28	12	12	0	0
5	A	158	0	0	3	2
5	B	171	0	0	3	3
All	All	3152	2773	2769	23	3

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 4.

All (23) 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:B:88:LYS:NZ	5:B:462:HOH:O	1.87	0.93
1:A:1:MET:SD	1:A:43[B]:GLN:NE2	2.49	0.85
1:A:107:GLU:OE1	5:A:456:HOH:O	1.98	0.82
1:A:73:ARG:NH2	5:A:386:HOH:O	2.20	0.74
1:B:94:HIS:CD2	5:B:429:HOH:O	2.44	0.71
1:A:45:VAL:HG22	1:A:50[B]:THR:HG22	1.73	0.70
1:B:1:MET:SD	1:B:43:GLN:NE2	2.66	0.64
1:A:1:MET:HE2	1:A:51:SER:HA	1.80	0.63
1:A:109:VAL:CG2	1:A:110:PRO:HD2	2.32	0.59
1:A:1:MET:HE3	1:A:50[A]:THR:HG23	1.84	0.59
1:A:1:MET:HE2	1:A:51:SER:CA	2.38	0.54
1:B:45:VAL:HG22	1:B:50[B]:THR:HG22	1.90	0.54
1:A:1:MET:CE	1:A:51:SER:CA	2.86	0.53
1:A:109:VAL:HG23	1:A:110:PRO:HD2	1.94	0.48
1:A:147:LYS:HE3	5:A:431:HOH:O	2.12	0.48
1:B:94:HIS:HD2	5:B:429:HOH:O	1.90	0.48
1:A:1:MET:CE	1:A:43[B]:GLN:HE22	2.27	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:109:VAL:HG22	1:A:110:PRO:HD2	1.97	0.47
1:A:1:MET:HE1	1:A:43[B]:GLN:NE2	2.30	0.47
1:A:105:ASP:O	1:A:106:SER:OG	2.31	0.43
1:A:1:MET:CE	1:A:51:SER:HA	2.45	0.43
1:A:1:MET:CE	1:A:43[B]:GLN:NE2	2.82	0.42
1:A:109:VAL:CG2	1:A:110:PRO:CD	2.99	0.40

All (3) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:A:445:HOH:O	5:B:440:HOH:O[1_455]	1.78	0.42
5:B:367:HOH:O	5:B:377:HOH:O[1_655]	2.06	0.14
5:A:401:HOH:O	5:B:462:HOH:O[1_455]	2.17	0.03

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	171/170 (101%)	168 (98%)	3 (2%)	0	100	100
1	B	170/170 (100%)	167 (98%)	3 (2%)	0	100	100
All	All	341/340 (100%)	335 (98%)	6 (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	152/150 (101%)	152 (100%)	0	100	100
1	B	151/150 (101%)	151 (100%)	0	100	100
All	All	303/300 (101%)	303 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	94	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 8 ligands modelled in this entry, 4 are monoatomic - leaving 4 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
3	20G	A	203	1	26,26,26	2.19	9 (34%)	34,34,34	1.64	8 (23%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	GDP	A	204	2	24,30,30	0.99	1 (4%)	31,47,47	1.89	5 (16%)
3	20G	B	203	1	26,26,26	1.97	7 (26%)	34,34,34	1.40	5 (14%)
4	GDP	B	204	2	24,30,30	1.57	3 (12%)	31,47,47	2.39	9 (29%)

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
3	20G	A	203	1	-	2/17/27/27	0/2/2/2
4	GDP	A	204	2	-	2/12/32/32	0/3/3/3
3	20G	B	203	1	-	2/17/27/27	0/2/2/2
4	GDP	B	204	2	-	3/12/32/32	0/3/3/3

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	203	20G	C15-N14	5.47	1.45	1.34
4	B	204	GDP	C2'-C1'	-5.37	1.45	1.53
3	A	203	20G	C09-N10	5.12	1.45	1.35
3	B	203	20G	C09-N10	4.62	1.44	1.35
3	B	203	20G	C03-C02	3.97	1.44	1.38
3	B	203	20G	C15-N14	3.73	1.42	1.34
3	A	203	20G	C16-C15	3.42	1.57	1.51
3	B	203	20G	C02-CL1	-3.15	1.67	1.74
3	B	203	20G	C04-CL2	-3.06	1.66	1.73
4	B	204	GDP	C6-C5	3.01	1.46	1.41
3	A	203	20G	C03-C02	2.73	1.42	1.38
3	A	203	20G	C03-C04	-2.66	1.34	1.38
4	A	204	GDP	O4'-C1'	2.53	1.44	1.41
3	A	203	20G	C04-CL2	-2.52	1.67	1.73
3	A	203	20G	C02-CL1	-2.52	1.69	1.74
3	B	203	20G	C13-N14	-2.47	1.41	1.46
3	A	203	20G	C18-S19	-2.38	1.72	1.80
3	B	203	20G	C22-N10	-2.35	1.42	1.47
4	B	204	GDP	C5-C4	2.19	1.46	1.40
3	A	203	20G	C12-C13	-2.02	1.47	1.52

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	204	GDP	C6-N1-C2	6.32	125.97	115.93
4	B	204	GDP	C5-C6-N1	-6.16	115.01	123.43
4	B	204	GDP	C6-C5-C4	-5.39	115.65	120.80
4	A	204	GDP	C5-C6-N1	-5.39	116.06	123.43
4	A	204	GDP	C6-N1-C2	5.25	124.27	115.93
4	A	204	GDP	C6-C5-C4	-4.42	116.57	120.80
4	B	204	GDP	C3'-C2'-C1'	4.19	107.28	100.98
4	B	204	GDP	C1'-N9-C4	-4.10	119.44	126.64
3	A	203	20G	C17-C16-C15	3.64	123.47	113.26
3	A	203	20G	C21-C22-N10	3.61	116.30	110.82
4	A	204	GDP	N3-C2-N1	-3.51	122.54	127.22
3	A	203	20G	C17-C18-S19	3.43	123.98	112.96
3	B	203	20G	C22-N10-C11	3.08	118.56	112.62
3	B	203	20G	C12-C13-N14	-2.72	104.93	110.56
4	B	204	GDP	O3'-C3'-C4'	2.66	118.75	111.05
3	B	203	20G	C17-C18-S19	2.63	121.41	112.96
3	B	203	20G	C24-C25-C02	2.62	122.01	119.24
3	A	203	20G	C16-C15-N14	2.61	120.36	115.83
3	A	203	20G	C24-C25-C02	2.49	121.87	119.24
4	B	204	GDP	O3'-C3'-C2'	-2.45	103.89	111.82
3	A	203	20G	C25-C24-C06	-2.29	115.68	120.06
4	A	204	GDP	C2-N3-C4	2.27	117.94	115.36
3	A	203	20G	O20-C15-C16	-2.17	118.04	122.02
4	B	204	GDP	N3-C2-N1	-2.16	124.34	127.22
3	A	203	20G	C18-C17-C16	2.16	120.13	113.38
3	B	203	20G	C25-C02-C03	-2.08	118.77	121.53
4	B	204	GDP	O2'-C2'-C3'	-2.07	105.12	111.82

There are no chirality outliers.

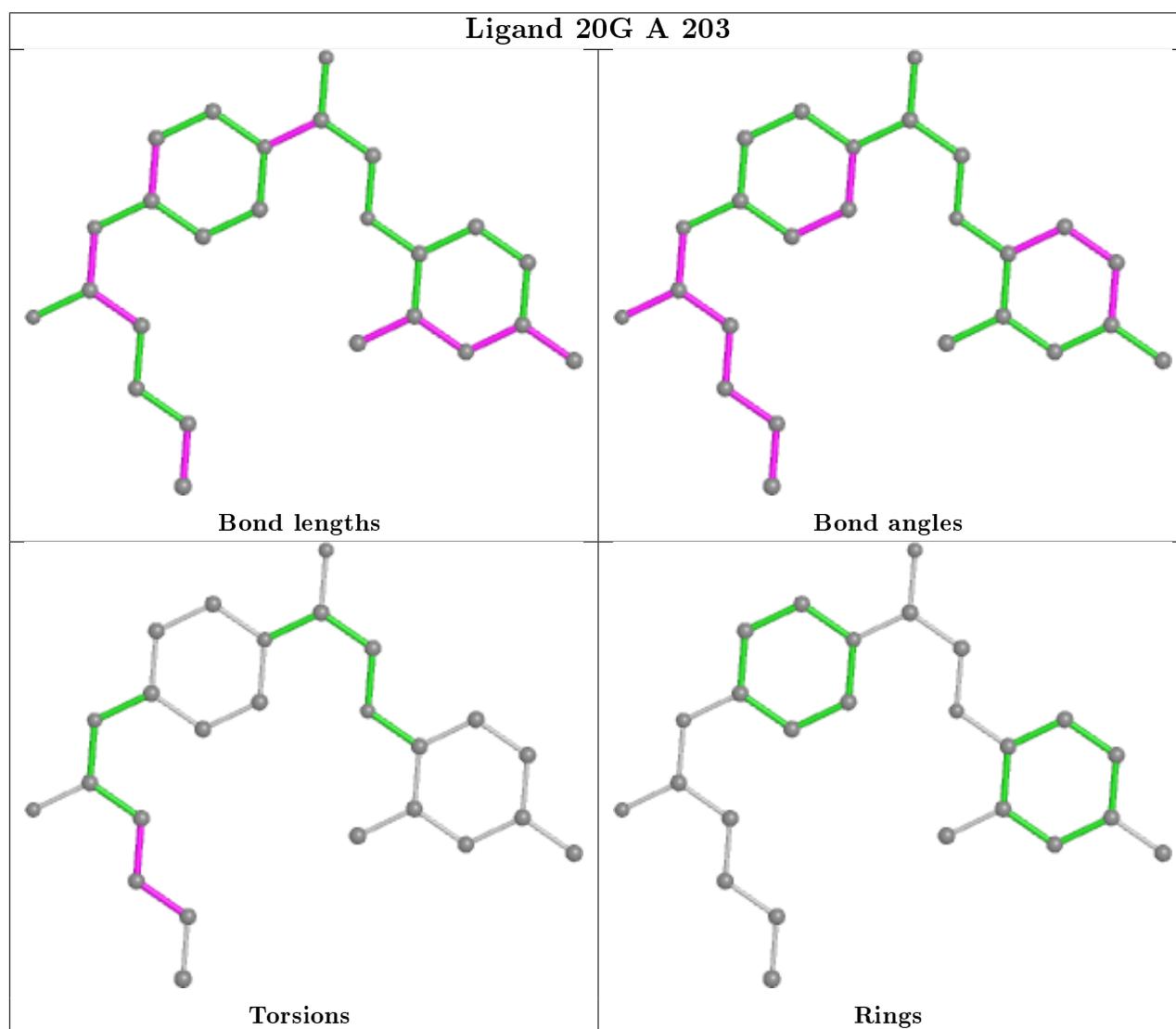
All (9) torsion outliers are listed below:

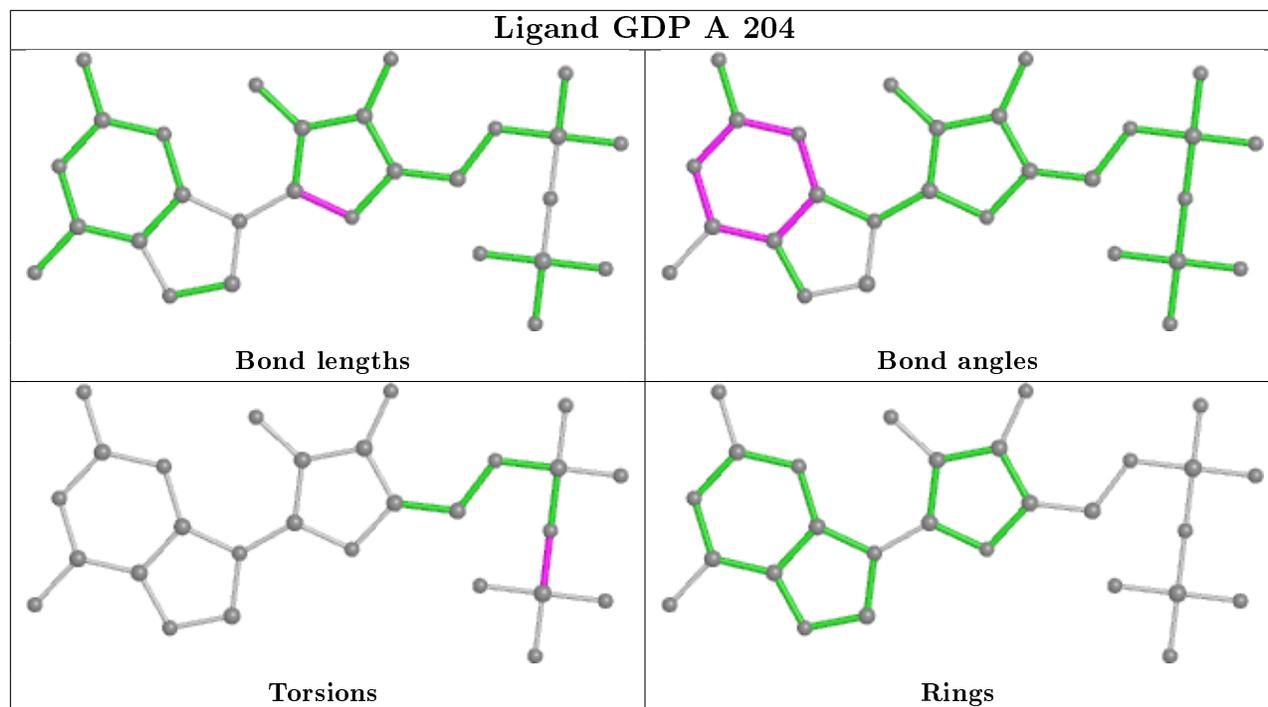
Mol	Chain	Res	Type	Atoms
3	A	203	20G	C15-C16-C17-C18
3	A	203	20G	C16-C17-C18-S19
4	A	204	GDP	PA-O3A-PB-O3B
3	B	203	20G	C15-C16-C17-C18
4	B	204	GDP	PA-O3A-PB-O1B
4	A	204	GDP	PA-O3A-PB-O1B
4	B	204	GDP	PA-O3A-PB-O2B
4	B	204	GDP	PA-O3A-PB-O3B
3	B	203	20G	C16-C17-C18-S19

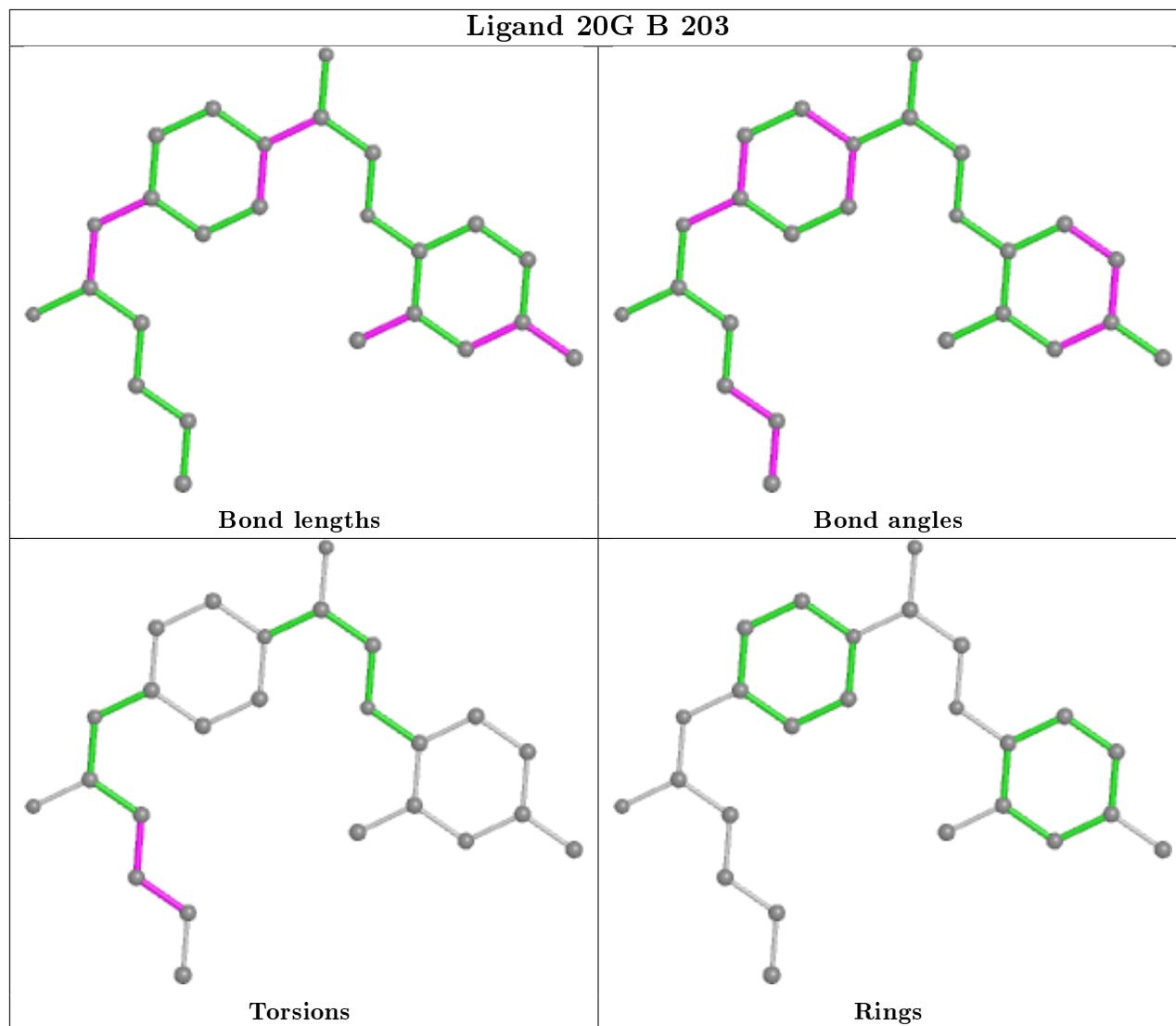
There are no ring outliers.

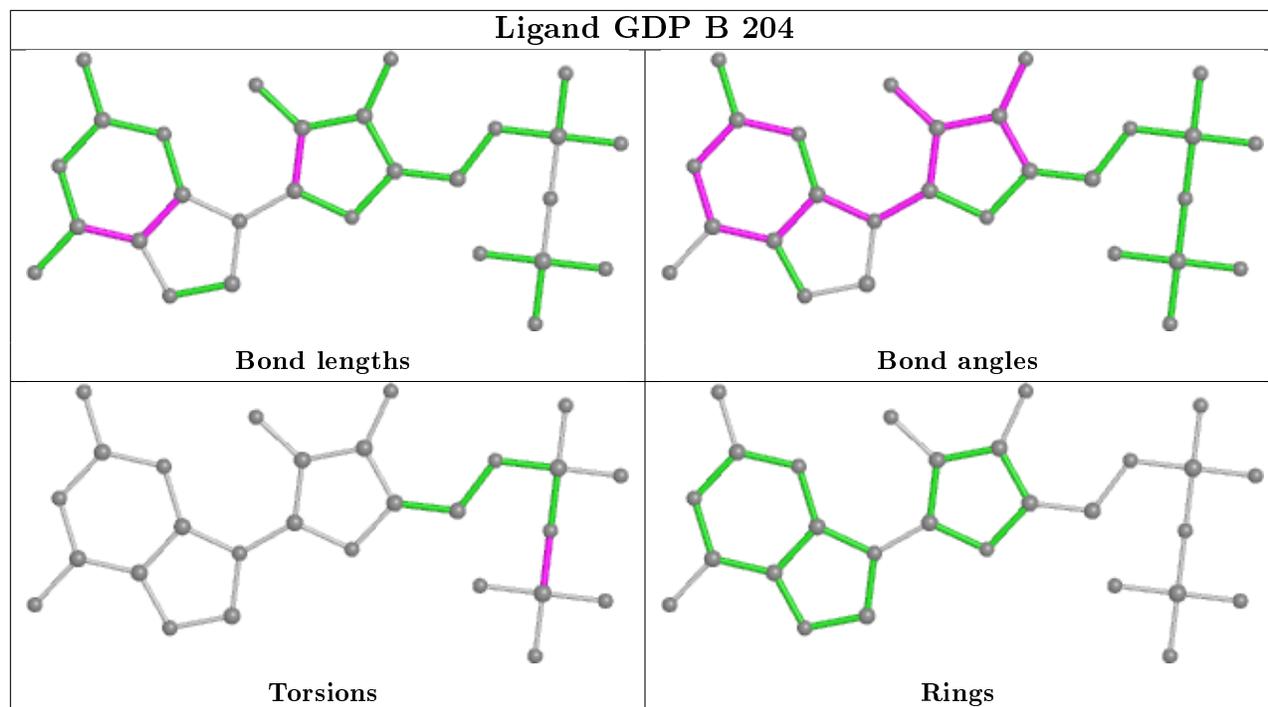
No monomer is involved in short contacts.

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.









5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

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

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

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	168/170 (98%)	0.02	2 (1%) 79 80	9, 16, 30, 46	0
1	B	168/170 (98%)	0.14	6 (3%) 42 39	10, 17, 36, 45	0
All	All	336/340 (98%)	0.08	8 (2%) 59 58	9, 16, 34, 46	0

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1	MET	5.0
1	B	1	MET	4.7
1	B	121	PRO	3.8
1	B	167	LYS	3.5
1	B	168	GLU	3.0
1	A	2	THR	2.9
1	B	122	SER	2.7
1	B	2	THR	2.3

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.

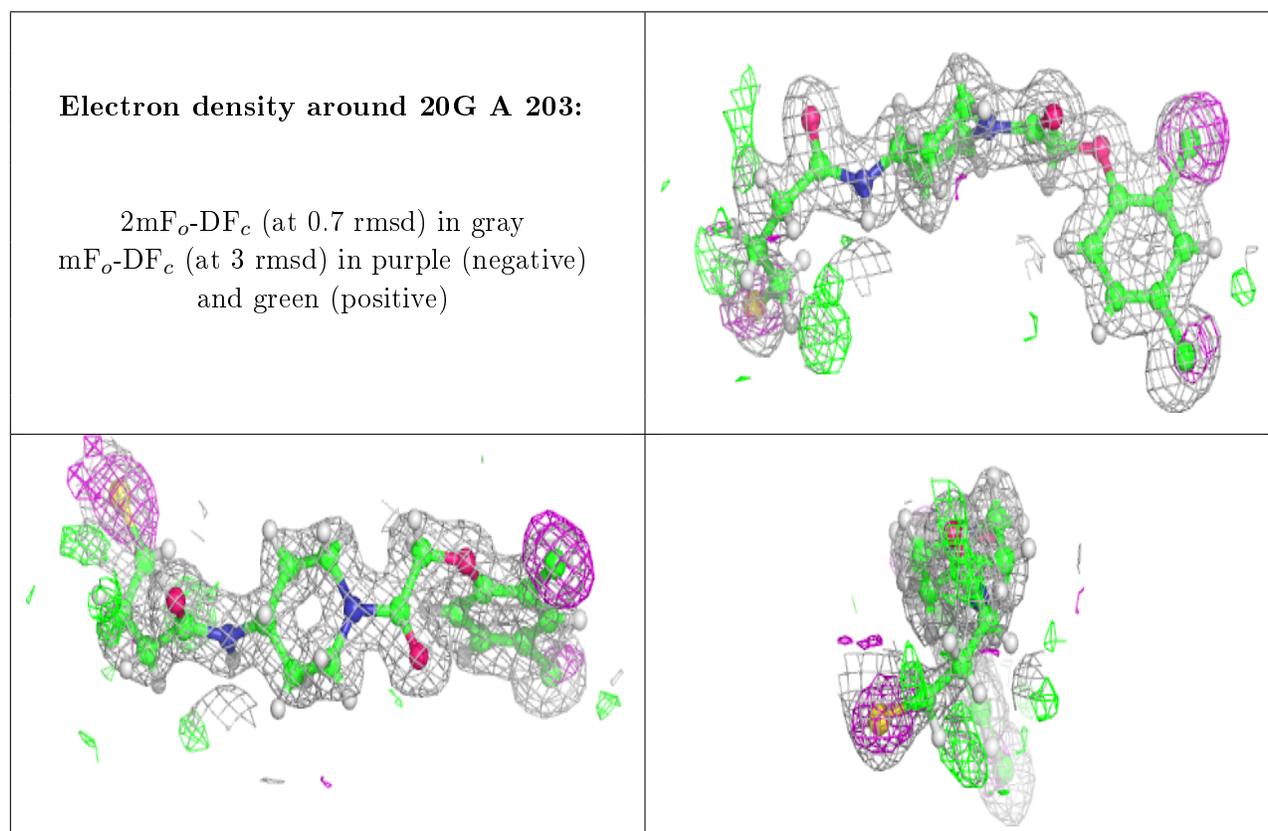
6.4 Ligands [i](#)

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.

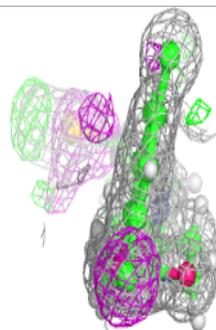
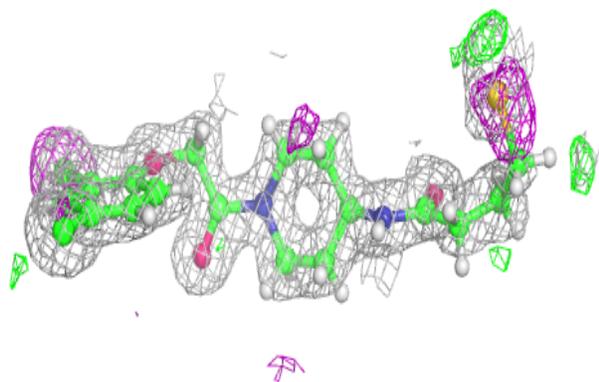
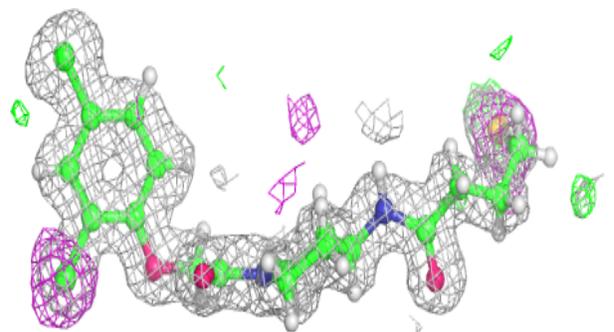
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	20G	A	203	25/25	0.94	0.09	14,18,29,31	0
3	20G	B	203	25/25	0.94	0.09	14,19,30,32	0
4	GDP	B	204	28/28	0.98	0.07	13,16,24,29	0
4	GDP	A	204	28/28	0.99	0.07	10,14,19,23	0
2	CA	A	201	1/1	1.00	0.03	13,13,13,13	0
2	CA	B	201	1/1	1.00	0.03	14,14,14,14	0
2	CA	B	202	1/1	1.00	0.04	14,14,14,14	0
2	CA	A	202	1/1	1.00	0.04	13,13,13,13	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.

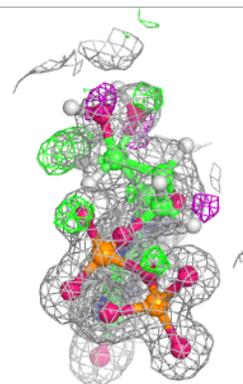
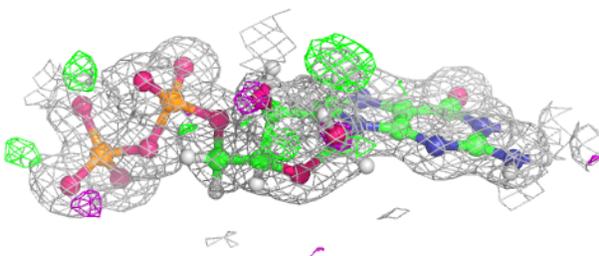
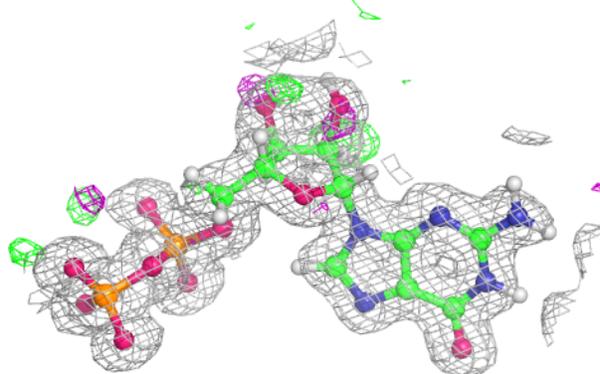


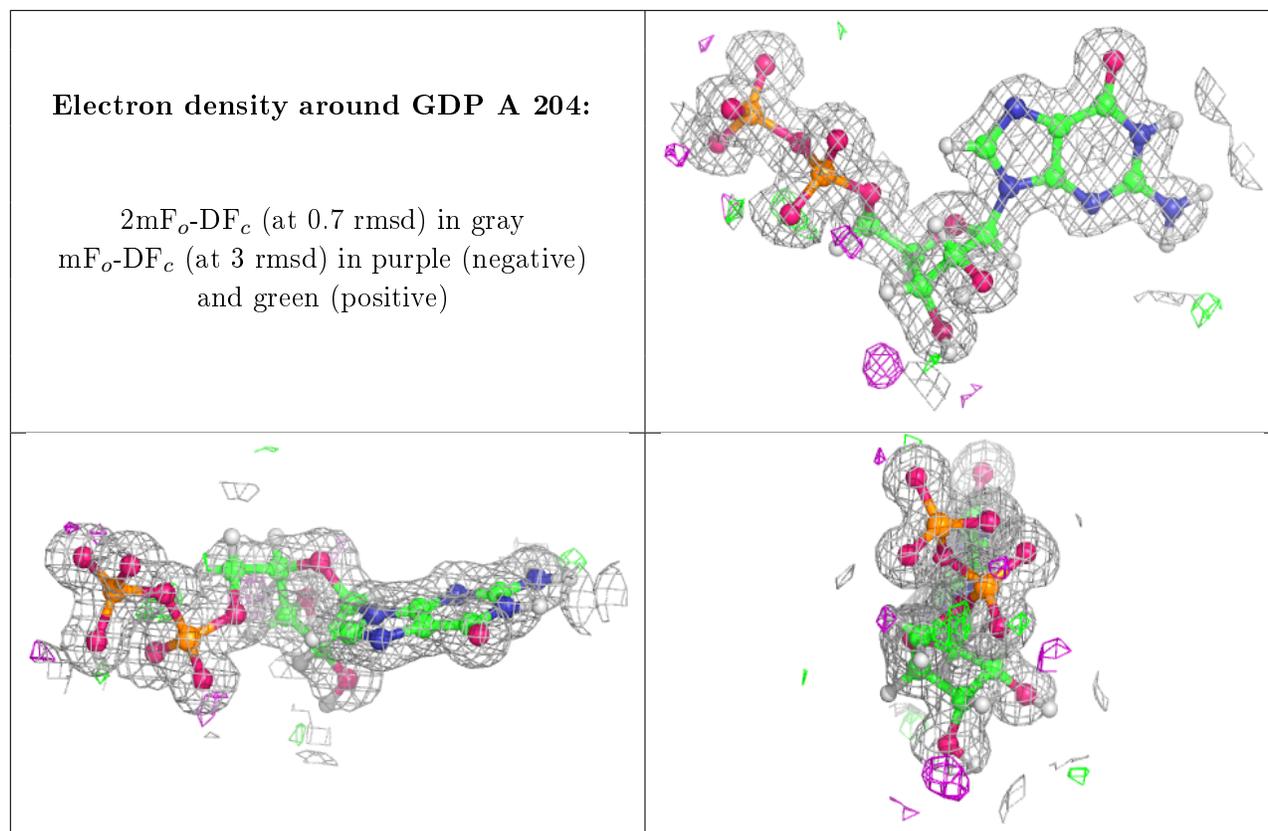
Electron density around 20G B 203:

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

**Electron density around GDP B 204:**

$2mF_o-DF_c$ (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.