



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 09:45 AM UTC

PDB ID : 4DFU / pdb_00004dfu
Title : Inhibition of an antibiotic resistance enzyme: crystal structure of aminoglycoside phosphotransferase APH(2")-ID/APH(2")-IVA in complex with kanamycin inhibited with quercetin
Authors : Stogios, P.J.; Minasov, G.; Dong, A.; Evdokimova, E.; Egorova, E.; Di Leo, R.; Li, H.; Shakya, T.; Wright, G.D.; Savchenko, A.; Anderson, W.F.; Center for Structural Genomics of Infectious Diseases (CSGID)
Deposited on : 2012-01-24
Resolution : 1.98 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)

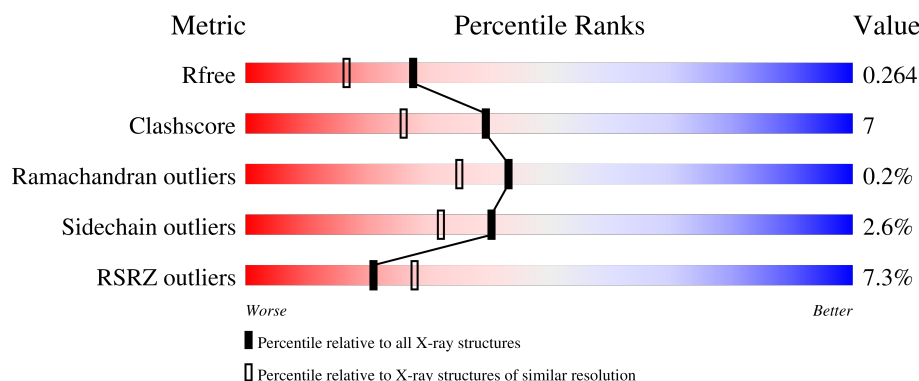
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.98 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	1506 (1.98-1.98)
Clashscore	190562	1534 (1.98-1.98)
Ramachandran outliers	187476	1518 (1.98-1.98)
Sidechain outliers	187428	1518 (1.98-1.98)
RSRZ outliers	180081	1506 (1.98-1.98)

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	322	7% 77% 14% 8%
1	B	322	6% 77% 15% 8%

Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 5268 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 APH(2")-Id.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	296	Total	C	N	O	S	Se	0	5	0
			2485	1603	400	470	3	9			
1	B	297	Total	C	N	O	S	Se	0	0	0
			2459	1587	396	466	3	7			

There are 42 discrepancies between the modelled and reference sequences:

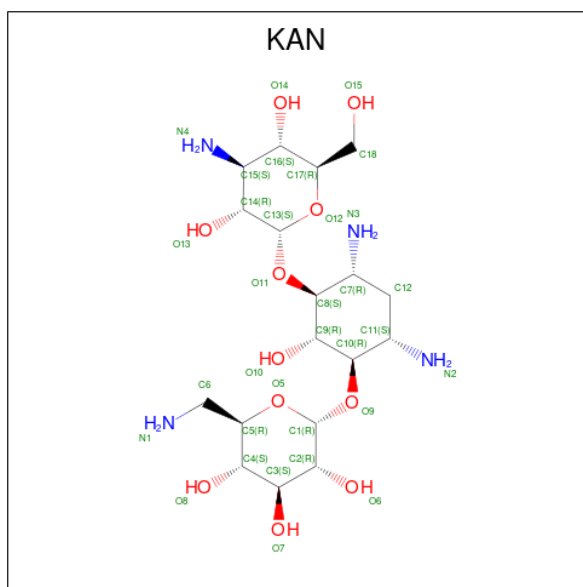
Chain	Residue	Modelled	Actual	Comment	Reference
A	-20	MSE	-	expression tag	UNP O68183
A	-19	GLY	-	expression tag	UNP O68183
A	-18	SER	-	expression tag	UNP O68183
A	-17	SER	-	expression tag	UNP O68183
A	-16	HIS	-	expression tag	UNP O68183
A	-15	HIS	-	expression tag	UNP O68183
A	-14	HIS	-	expression tag	UNP O68183
A	-13	HIS	-	expression tag	UNP O68183
A	-12	HIS	-	expression tag	UNP O68183
A	-11	HIS	-	expression tag	UNP O68183
A	-10	SER	-	expression tag	UNP O68183
A	-9	SER	-	expression tag	UNP O68183
A	-8	GLY	-	expression tag	UNP O68183
A	-7	ARG	-	expression tag	UNP O68183
A	-6	GLU	-	expression tag	UNP O68183
A	-5	ASN	-	expression tag	UNP O68183
A	-4	LEU	-	expression tag	UNP O68183
A	-3	TYR	-	expression tag	UNP O68183
A	-2	PHE	-	expression tag	UNP O68183
A	-1	GLN	-	expression tag	UNP O68183
A	0	GLY	-	expression tag	UNP O68183
B	-20	MSE	-	expression tag	UNP O68183
B	-19	GLY	-	expression tag	UNP O68183
B	-18	SER	-	expression tag	UNP O68183
B	-17	SER	-	expression tag	UNP O68183

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-16	HIS	-	expression tag	UNP O68183
B	-15	HIS	-	expression tag	UNP O68183
B	-14	HIS	-	expression tag	UNP O68183
B	-13	HIS	-	expression tag	UNP O68183
B	-12	HIS	-	expression tag	UNP O68183
B	-11	HIS	-	expression tag	UNP O68183
B	-10	SER	-	expression tag	UNP O68183
B	-9	SER	-	expression tag	UNP O68183
B	-8	GLY	-	expression tag	UNP O68183
B	-7	ARG	-	expression tag	UNP O68183
B	-6	GLU	-	expression tag	UNP O68183
B	-5	ASN	-	expression tag	UNP O68183
B	-4	LEU	-	expression tag	UNP O68183
B	-3	TYR	-	expression tag	UNP O68183
B	-2	PHE	-	expression tag	UNP O68183
B	-1	GLN	-	expression tag	UNP O68183
B	0	GLY	-	expression tag	UNP O68183

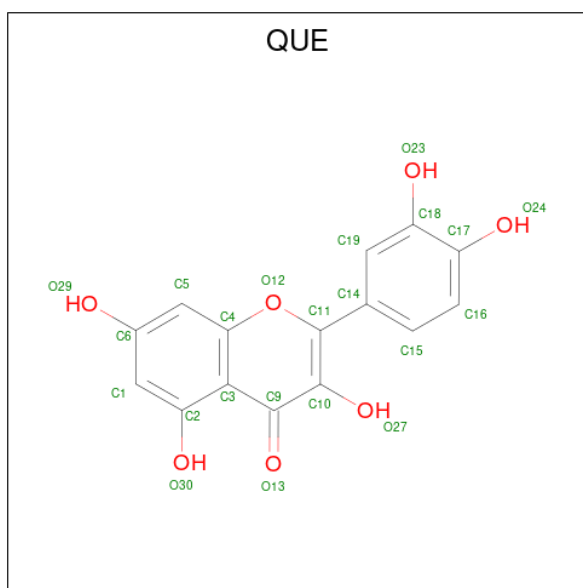
- Molecule 2 is KANAMYCIN A (CCD ID: KAN) (formula: $C_{18}H_{36}N_4O_{11}$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			33	18	4	11		
2	B	1	Total	C	N	O	0	0
			33	18	4	11		

- Molecule 3 is 3,5,7,3',4'-PENTAHYDROXYFLAVONE (CCD ID: QUE) (formula:

C₁₅H₁₀O₇).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 22 15 7	0	0
3	B	1	Total C O 44 30 14	0	1
3	B	1	Total C O 22 15 7	0	0

- Molecule 4 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total Cl 1 1	0	0
4	B	1	Total Cl 1 1	0	0

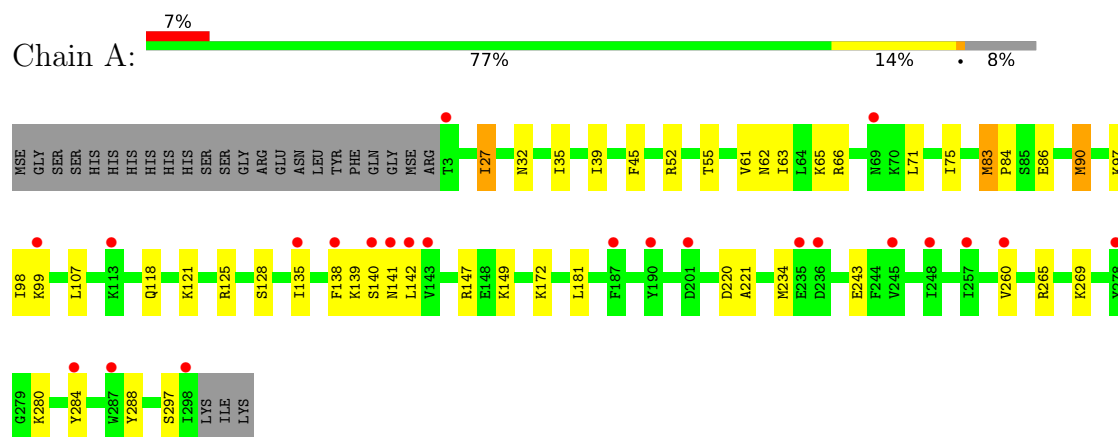
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	74	Total O 74 74	0	0
5	B	94	Total O 94 94	0	0

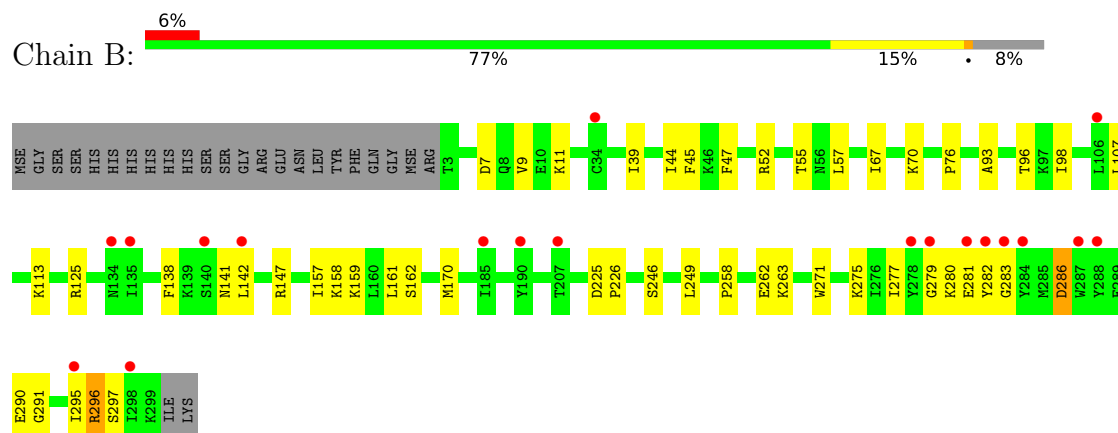
3 Residue-property plots

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

• Molecule 1: APH(2")-Id



• Molecule 1: APH(2")-Id



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	43.14Å 101.81Å 70.25Å 90.00° 96.86° 90.00°	Depositor
Resolution (Å)	19.84 – 1.98 19.84 – 1.98	Depositor EDS
% Data completeness (in resolution range)	98.8 (19.84-1.98) 90.9 (19.84-1.98)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.04	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.60 (at 1.97Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7.3_928)	Depositor
R, R_{free}	0.219 , 0.270 0.214 , 0.264	Depositor DCC
R_{free} test set	1995 reflections (4.77%)	wwPDB-VP
Wilson B-factor (Å ²)	35.3	Xtriage
Anisotropy	0.460	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 59.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	5268	wwPDB-VP
Average B, all atoms (Å ²)	68.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.71% 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: QUE, KAN, CL

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	0.45	1/2543 (0.0%)	0.78	0/3416
1	B	0.47	0/2511	0.76	0/3373
All	All	0.46	1/5054 (0.0%)	0.77	0/6789

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	90	MSE	SE-CE	-5.61	1.78	1.95

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	2485	0	2445	33	0
1	B	2459	0	2417	34	0
2	A	33	0	36	1	0
2	B	33	0	36	0	0
3	A	22	0	7	2	0
3	B	66	0	19	13	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	A	74	0	0	1	0
5	B	94	0	0	1	0
All	All	5268	0	4960	67	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 7.

All (67) 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:140:SER:HB3	1:B:141:ASN:HA	1.63	0.81
1:A:62[B]:ASN:OD1	1:A:66:ARG:NH1	2.15	0.80
1:A:98:ILE:N	3:A:402:QUE:O13	2.16	0.78
1:A:98:ILE:O	3:A:402:QUE:O27	2.07	0.70
1:A:147:ARG:NH2	1:A:181:LEU:O	2.26	0.68
1:B:138:PHE:HB3	1:B:142:LEU:HD12	1.75	0.67
1:A:172:LYS:NZ	1:A:297:SER:O	2.28	0.66
1:A:234:MSE:HE1	1:A:260:VAL:HG12	1.79	0.64
1:B:52:ARG:NE	1:B:281:GLU:OE1	2.32	0.63
1:A:32:ASN:ND2	1:A:284:TYR:OH	2.31	0.62
1:B:67:ILE:HA	1:B:70:LYS:HD2	1.83	0.60
1:B:161:LEU:HB2	1:B:170:MSE:HE1	1.86	0.58
1:A:62[B]:ASN:HD21	1:A:142:LEU:HD11	1.70	0.55
1:A:63:ILE:HD13	1:A:221:ALA:HB3	1.88	0.54
1:B:7:ASP:O	1:B:11:LYS:HG3	2.07	0.54
1:B:162:SER:HA	1:B:170:MSE:HE3	1.89	0.53
1:A:66:ARG:HH22	1:A:139:LYS:HG3	1.74	0.53
1:B:281:GLU:HG2	1:B:282:TYR:CD2	2.45	0.52
1:B:258:PRO:O	1:B:262:GLU:HG3	2.09	0.52
1:B:98:ILE:N	3:B:403:QUE:O13	2.21	0.51
1:B:225:ASP:OD2	1:B:263:LYS:NZ	2.44	0.50
1:B:271:TRP:NE1	1:B:275:LYS:HD2	2.27	0.50
1:A:220:ASP:OD2	2:A:401:KAN:N4	2.40	0.50
1:A:97:LYS:HD3	1:A:99:LYS:HZ3	1.77	0.49
1:A:55:THR:HG21	3:B:401[B]:QUE:C9	2.42	0.49
1:A:83[A]:MSE:HE2	1:A:84:PRO:HD3	1.93	0.49
1:B:39:ILE:HD12	1:B:45:PHE:CD1	2.48	0.49
1:B:55:THR:HG21	3:B:401[A]:QUE:C9	2.42	0.48
1:A:86:GLU:H	1:A:86:GLU:CD	2.21	0.48
1:A:61:VAL:HG21	1:A:90:MSE:CE	2.44	0.47
1:A:118:GLN:NE2	1:A:121:LYS:HD3	2.29	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:403:QUE:O27	3:B:403:QUE:H19	2.14	0.47
1:A:265:ARG:HG2	1:A:269:LYS:HE3	1.97	0.47
1:B:246:SER:HA	1:B:249:LEU:HD12	1.96	0.47
1:A:141:ASN:ND2	1:B:141:ASN:HB3	2.30	0.46
1:B:76:PRO:HB3	3:B:403:QUE:H1	1.97	0.46
1:B:296:ARG:HG3	1:B:297:SER:N	2.31	0.46
1:A:27:ILE:HG12	1:A:35:ILE:HD11	1.98	0.46
1:A:83[A]:MSE:HE3	1:B:159:LYS:HD3	1.98	0.45
1:A:135:ILE:O	1:A:138:PHE:HB2	2.17	0.45
1:A:65[A]:LYS:NZ	5:A:529:HOH:O	2.50	0.45
1:B:158:LYS:O	1:B:162:SER:HB3	2.17	0.44
1:B:286:ASP:O	1:B:290:GLU:HG3	2.17	0.44
3:B:401[B]:QUE:H15	3:B:401[B]:QUE:O27	2.18	0.44
1:A:280:LYS:HG2	1:A:288:TYR:CD1	2.53	0.44
1:B:96:THR:O	3:B:403:QUE:O30	2.35	0.44
1:A:55:THR:HG21	3:B:401[B]:QUE:C10	2.48	0.43
1:A:62[B]:ASN:CG	1:A:139:LYS:HD2	2.44	0.43
1:A:141:ASN:CG	1:B:141:ASN:HB3	2.43	0.43
1:B:9:VAL:HG13	1:B:47:PHE:CE1	2.54	0.43
1:A:39:ILE:HD12	1:A:45:PHE:CD1	2.54	0.42
1:A:66:ARG:NH2	1:A:139:LYS:HG3	2.33	0.42
1:A:118:GLN:HE22	1:A:121:LYS:HD3	1.83	0.42
1:A:71:LEU:HD12	1:A:75:ILE:HD13	2.01	0.42
1:B:125:ARG:HD2	5:B:575:HOH:O	2.19	0.42
3:B:401[B]:QUE:O13	3:B:401[B]:QUE:O30	2.36	0.41
1:B:98:ILE:O	3:B:403:QUE:O27	2.37	0.41
1:B:281:GLU:C	1:B:283:GLY:H	2.28	0.41
1:B:52:ARG:NE	3:B:401[B]:QUE:O24	2.53	0.41
1:B:296:ARG:HG3	1:B:297:SER:H	1.84	0.41
1:A:55:THR:HG21	3:B:401[A]:QUE:C11	2.50	0.41
1:B:291:GLY:O	1:B:295:ILE:HG12	2.20	0.41
3:B:403:QUE:O13	3:B:403:QUE:O30	2.38	0.41
1:B:57:LEU:HD11	1:B:93:ALA:HB3	2.04	0.40
1:B:157:ILE:HG12	1:B:277:ILE:HD11	2.02	0.40
1:B:113:LYS:HE3	1:B:113:LYS:HB2	1.76	0.40
1:B:225:ASP:HA	1:B:226:PRO:HD2	1.92	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	299/322 (93%)	285 (95%)	14 (5%)	0	100	100
1	B	295/322 (92%)	280 (95%)	14 (5%)	1 (0%)	36	27
All	All	594/644 (92%)	565 (95%)	28 (5%)	1 (0%)	43	35

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	279	GLY

5.3.2 Protein sidechains

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	279/288 (97%)	270 (97%)	9 (3%)	34	24
1	B	275/288 (96%)	269 (98%)	6 (2%)	45	38
All	All	554/576 (96%)	539 (97%)	15 (3%)	40	31

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	27	ILE
1	A	52	ARG
1	A	83[A]	MSE
1	A	83[B]	MSE
1	A	107	LEU

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Mol	Chain	Res	Type
1	A	125	ARG
1	A	128	SER
1	A	149	LYS
1	A	243	GLU
1	B	44	ILE
1	B	107	LEU
1	B	147	ARG
1	B	280	LYS
1	B	286	ASP
1	B	296	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	32	ASN
1	A	141	ASN
1	A	250	ASN
1	B	117	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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 8 ligands modelled in this entry, 2 are monoatomic - leaving 6 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
2	KAN	B	402	-	34,35,35	0.71	0	47,52,52	1.36	9 (19%)
3	QUE	B	403	-	24,24,24	1.76	6 (25%)	36,36,36	2.97	8 (22%)
3	QUE	B	401[B]	-	24,24,24	1.63	5 (20%)	36,36,36	2.88	9 (25%)
3	QUE	A	402	-	24,24,24	1.70	6 (25%)	36,36,36	2.94	10 (27%)
3	QUE	B	401[A]	-	24,24,24	1.72	4 (16%)	36,36,36	2.79	9 (25%)
2	KAN	A	401	-	34,35,35	0.64	0	47,52,52	1.24	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
2	KAN	B	402	-	-	4/12/72/72	0/3/3/3
3	QUE	B	403	-	-	0/4/4/4	0/3/3/3
3	QUE	B	401[B]	-	-	0/4/4/4	0/3/3/3
3	QUE	A	402	-	-	0/4/4/4	0/3/3/3
3	QUE	B	401[A]	-	-	0/4/4/4	0/3/3/3
2	KAN	A	401	-	-	4/12/72/72	0/3/3/3

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	403	QUE	C14-C11	-5.52	1.39	1.47
3	B	401[A]	QUE	C14-C11	-5.27	1.39	1.47
3	A	402	QUE	C14-C11	-5.08	1.39	1.47
3	B	401[B]	QUE	C14-C11	-5.01	1.39	1.47
3	A	402	QUE	O12-C4	-3.70	1.32	1.38
3	B	401[A]	QUE	O12-C4	-3.61	1.33	1.38
3	B	401[A]	QUE	C10-C11	3.60	1.40	1.36
3	B	403	QUE	O12-C4	-3.49	1.33	1.38
3	B	401[B]	QUE	O12-C4	-3.43	1.33	1.38
3	B	401[B]	QUE	C10-C11	3.35	1.39	1.36
3	A	402	QUE	C10-C11	2.90	1.39	1.36
3	B	403	QUE	C3-C9	-2.72	1.39	1.46
3	B	403	QUE	C10-C9	-2.54	1.39	1.46
3	B	403	QUE	C10-C11	2.48	1.38	1.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	401[B]	QUE	C3-C9	-2.43	1.40	1.46
3	B	401[A]	QUE	C3-C9	-2.43	1.40	1.46
3	A	402	QUE	O12-C11	-2.38	1.34	1.37
3	A	402	QUE	C10-C9	-2.31	1.40	1.46
3	A	402	QUE	C3-C9	-2.23	1.40	1.46
3	B	403	QUE	O12-C11	-2.13	1.35	1.37
3	B	401[B]	QUE	C10-C9	-2.11	1.40	1.46

All (53) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	403	QUE	O12-C4-C3	-10.33	111.06	121.19
3	A	402	QUE	O12-C4-C3	-9.88	111.51	121.19
3	B	401[B]	QUE	O12-C4-C3	-9.87	111.51	121.19
3	B	401[A]	QUE	O12-C4-C3	-9.40	111.98	121.19
3	A	402	QUE	C11-C10-C9	-7.68	116.03	121.52
3	B	403	QUE	C11-C10-C9	-7.31	116.30	121.52
3	B	403	QUE	O12-C4-C5	7.27	126.13	115.83
3	B	401[A]	QUE	C11-C10-C9	-7.18	116.39	121.52
3	B	401[B]	QUE	O12-C4-C5	6.99	125.73	115.83
3	B	401[B]	QUE	C11-C10-C9	-6.87	116.61	121.52
3	A	402	QUE	O12-C4-C5	6.55	125.11	115.83
3	B	401[A]	QUE	O12-C4-C5	6.49	125.03	115.83
3	A	402	QUE	O12-C11-C10	-6.16	114.35	120.19
3	A	402	QUE	C4-O12-C11	5.40	128.94	120.67
3	B	403	QUE	O12-C11-C10	-5.35	115.12	120.19
3	B	401[B]	QUE	O12-C11-C10	-5.35	115.12	120.19
3	B	401[A]	QUE	O12-C11-C10	-5.19	115.27	120.19
3	B	401[B]	QUE	C4-O12-C11	4.54	127.63	120.67
3	B	403	QUE	C4-O12-C11	4.42	127.44	120.67
3	B	403	QUE	O27-C10-C11	4.36	125.25	120.74
3	B	401[A]	QUE	C4-O12-C11	4.32	127.29	120.67
3	A	402	QUE	C3-C9-C10	3.80	122.27	116.88
3	B	401[B]	QUE	O12-C11-C14	3.61	115.93	110.87
3	B	401[B]	QUE	O27-C10-C11	3.45	124.31	120.74
2	B	402	KAN	O11-C8-C9	3.41	115.89	107.23
3	B	401[A]	QUE	O27-C10-C11	3.30	124.16	120.74
2	B	402	KAN	O9-C10-C9	3.19	115.34	107.23
3	B	403	QUE	O12-C11-C14	3.15	115.28	110.87
3	B	403	QUE	C3-C9-C10	3.06	121.21	116.88
3	B	401[B]	QUE	C3-C9-C10	3.03	121.18	116.88
3	B	401[A]	QUE	C3-C9-C10	2.99	121.11	116.88

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	401[A]	QUE	C19-C14-C11	-2.97	116.46	120.65
2	A	401	KAN	O9-C10-C9	2.92	114.65	107.23
2	A	401	KAN	O11-C8-C9	2.91	114.62	107.23
2	A	401	KAN	C17-C16-C15	-2.86	107.51	110.67
2	B	402	KAN	C6-C5-C4	-2.78	106.31	112.80
2	B	402	KAN	C1-O9-C10	2.65	124.26	117.98
3	A	402	QUE	O27-C10-C11	2.44	123.27	120.74
3	A	402	QUE	O12-C11-C14	2.41	114.25	110.87
2	A	401	KAN	C1-O9-C10	2.39	123.64	117.98
2	B	402	KAN	C10-C9-C8	2.33	113.84	109.11
2	B	402	KAN	C16-C15-C14	-2.28	106.11	111.06
2	A	401	KAN	C18-C17-C16	-2.25	107.49	113.02
2	B	402	KAN	O5-C5-C6	2.22	110.34	106.07
3	B	401[A]	QUE	O12-C11-C14	2.22	113.98	110.87
2	A	401	KAN	C4-C3-C2	-2.16	107.04	110.83
2	A	401	KAN	C16-C15-C14	-2.10	106.49	111.06
3	A	402	QUE	O13-C9-C3	-2.06	118.33	122.06
2	B	402	KAN	C11-C12-C7	2.05	114.54	111.02
2	A	401	KAN	C13-C14-C15	2.05	113.15	110.40
3	B	401[B]	QUE	C19-C14-C11	-2.04	117.77	120.65
3	A	402	QUE	O27-C10-C9	2.03	120.62	116.89
2	B	402	KAN	C18-C17-C16	-2.03	108.04	113.02

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	401	KAN	C9-C10-O9-C1
2	B	402	KAN	C9-C10-O9-C1
2	A	401	KAN	C11-C10-O9-C1
2	B	402	KAN	C11-C10-O9-C1
2	B	402	KAN	C2-C1-O9-C10
2	B	402	KAN	O5-C1-O9-C10
2	A	401	KAN	C2-C1-O9-C10
2	A	401	KAN	O5-C1-O9-C10

There are no ring outliers.

5 monomers are involved in 16 short contacts:

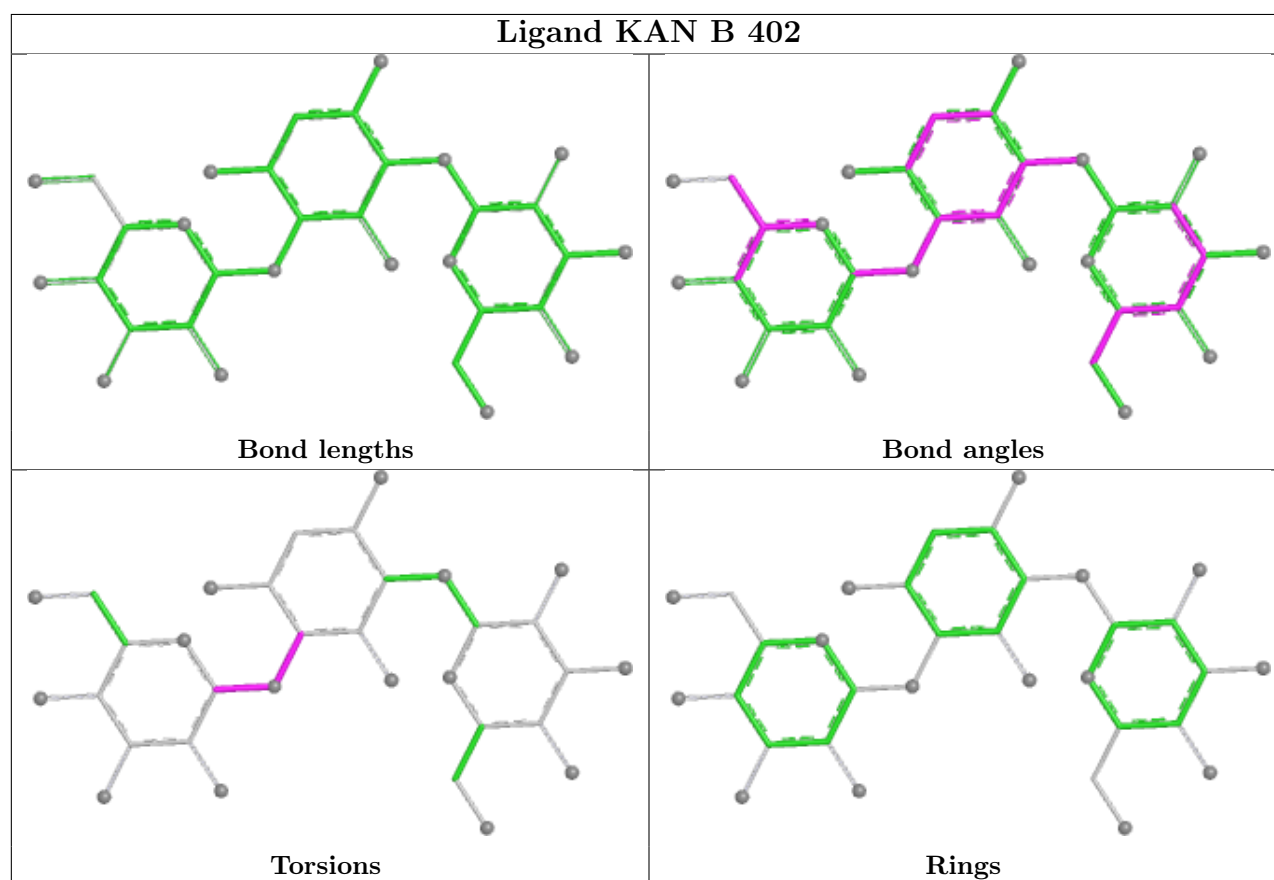
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	403	QUE	6	0

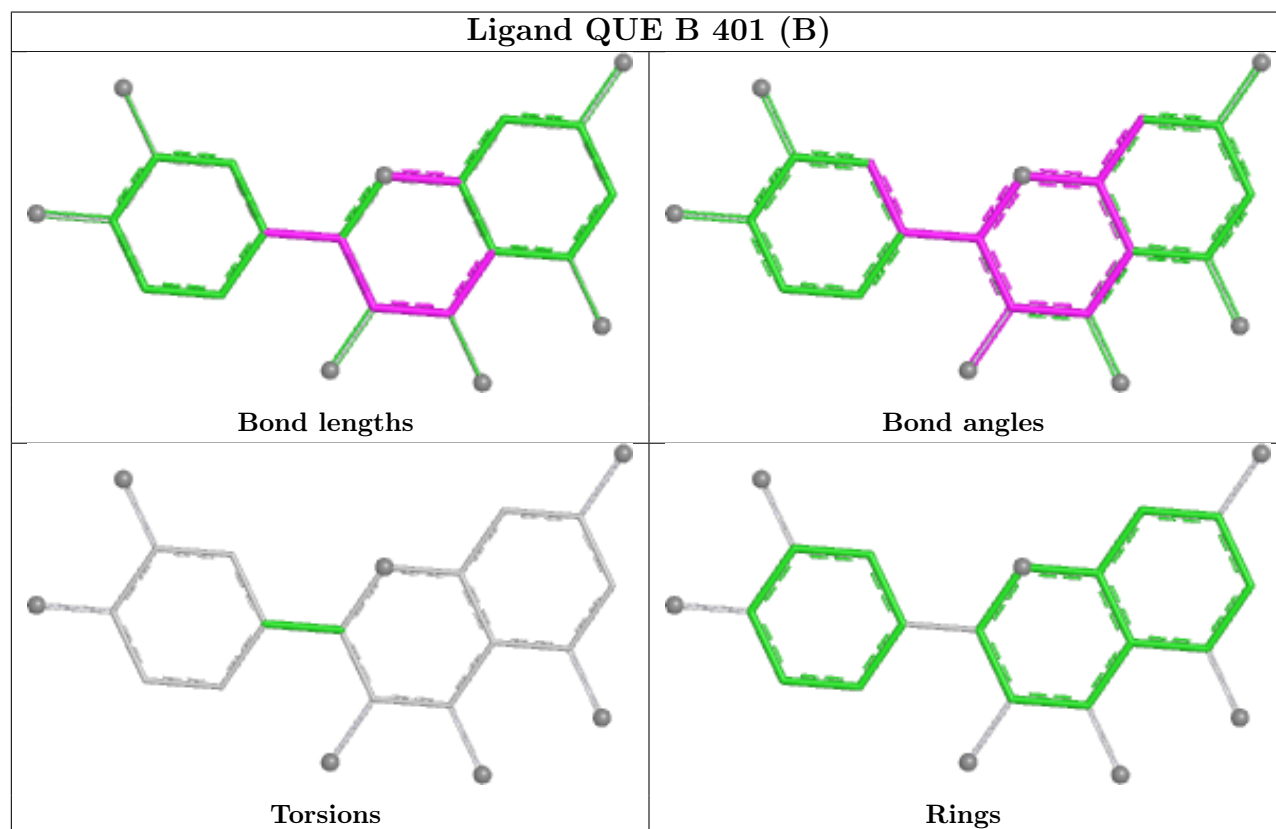
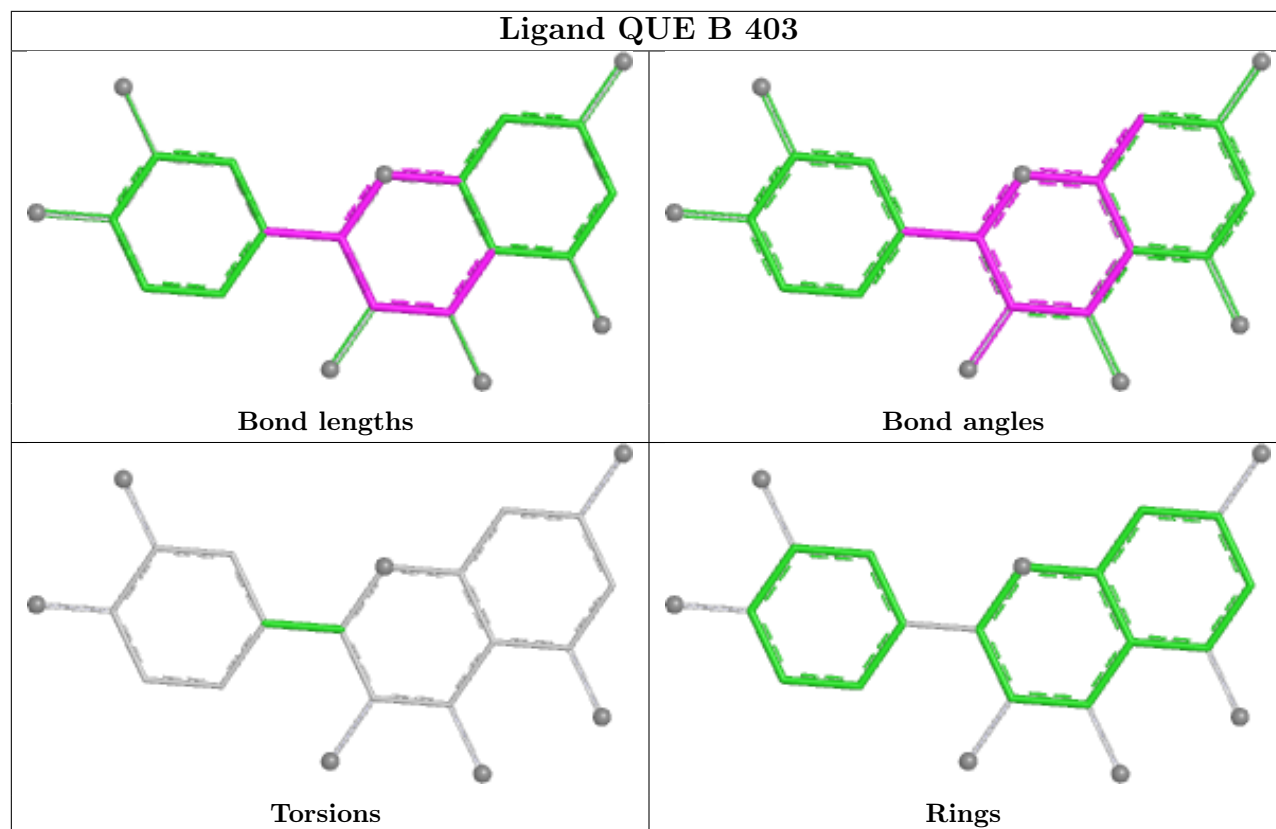
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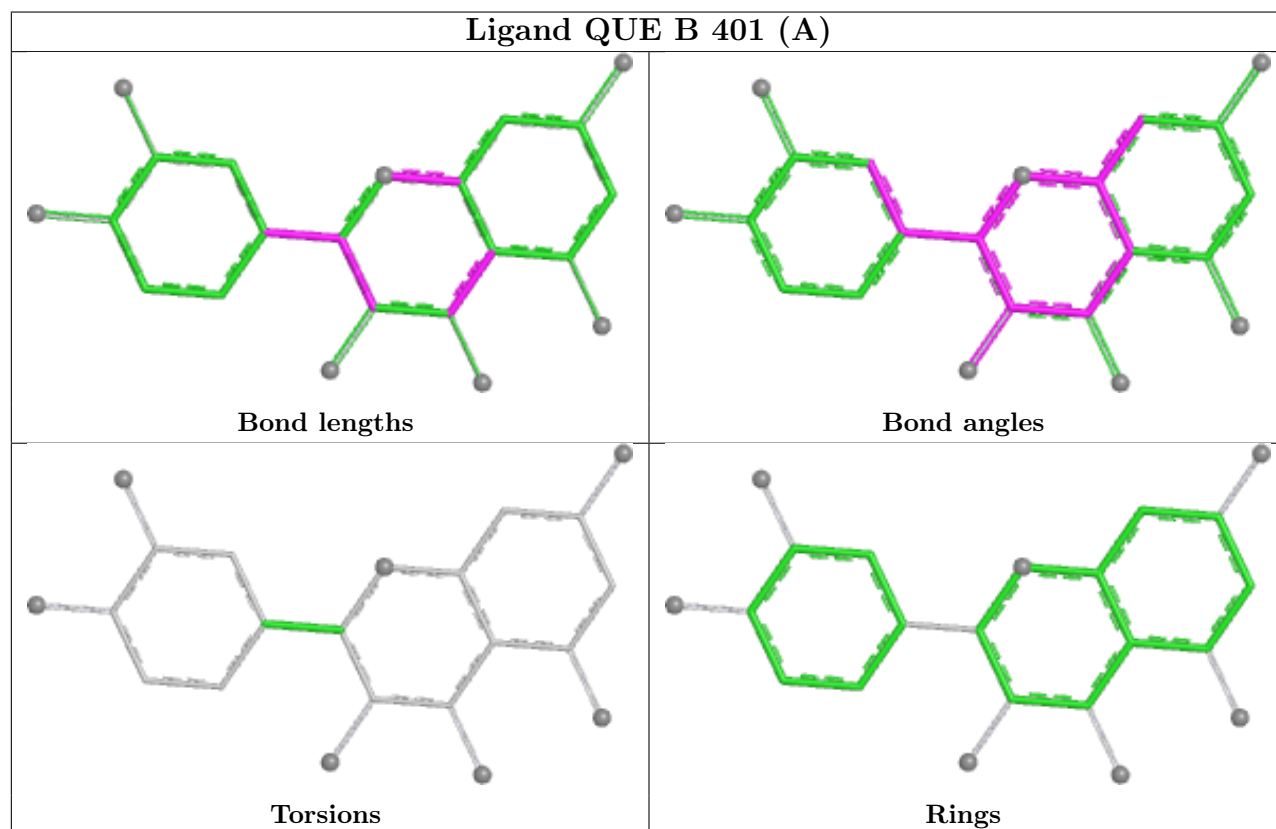
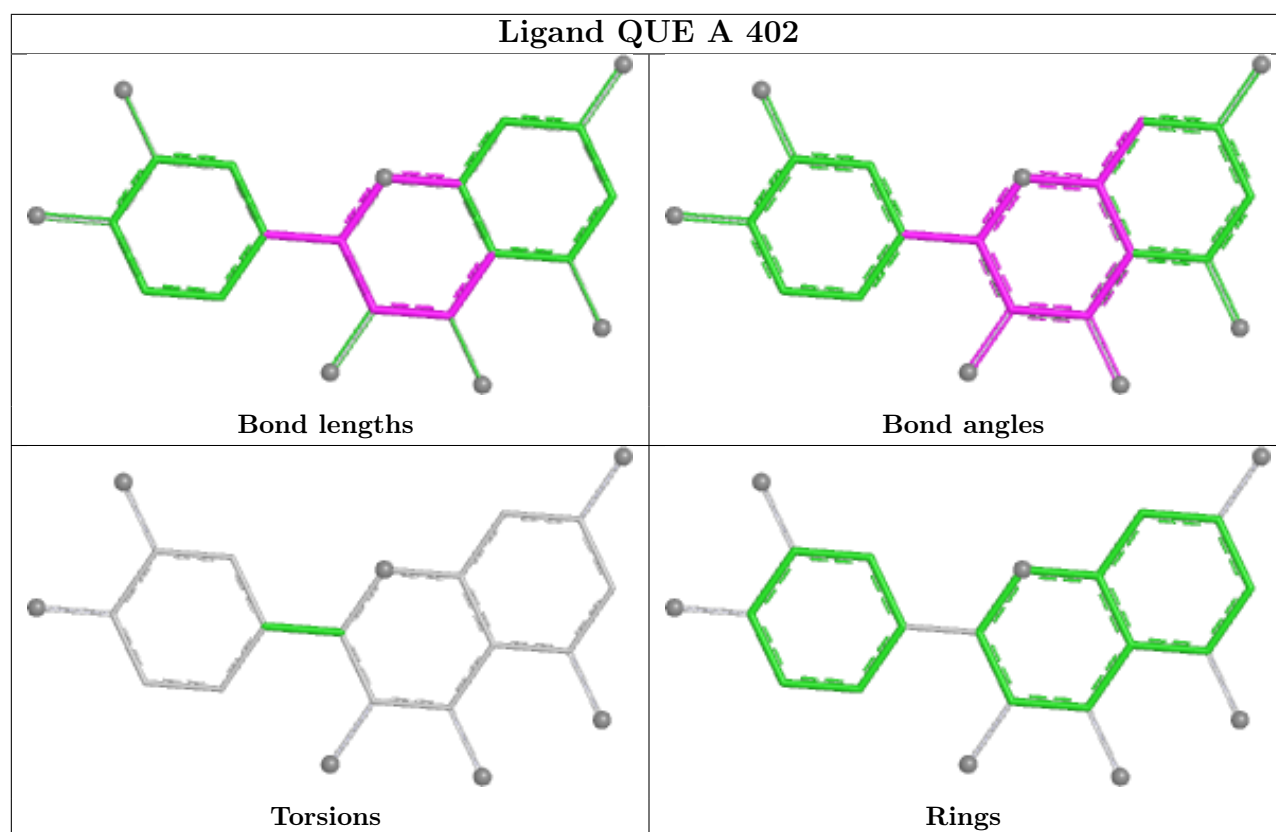
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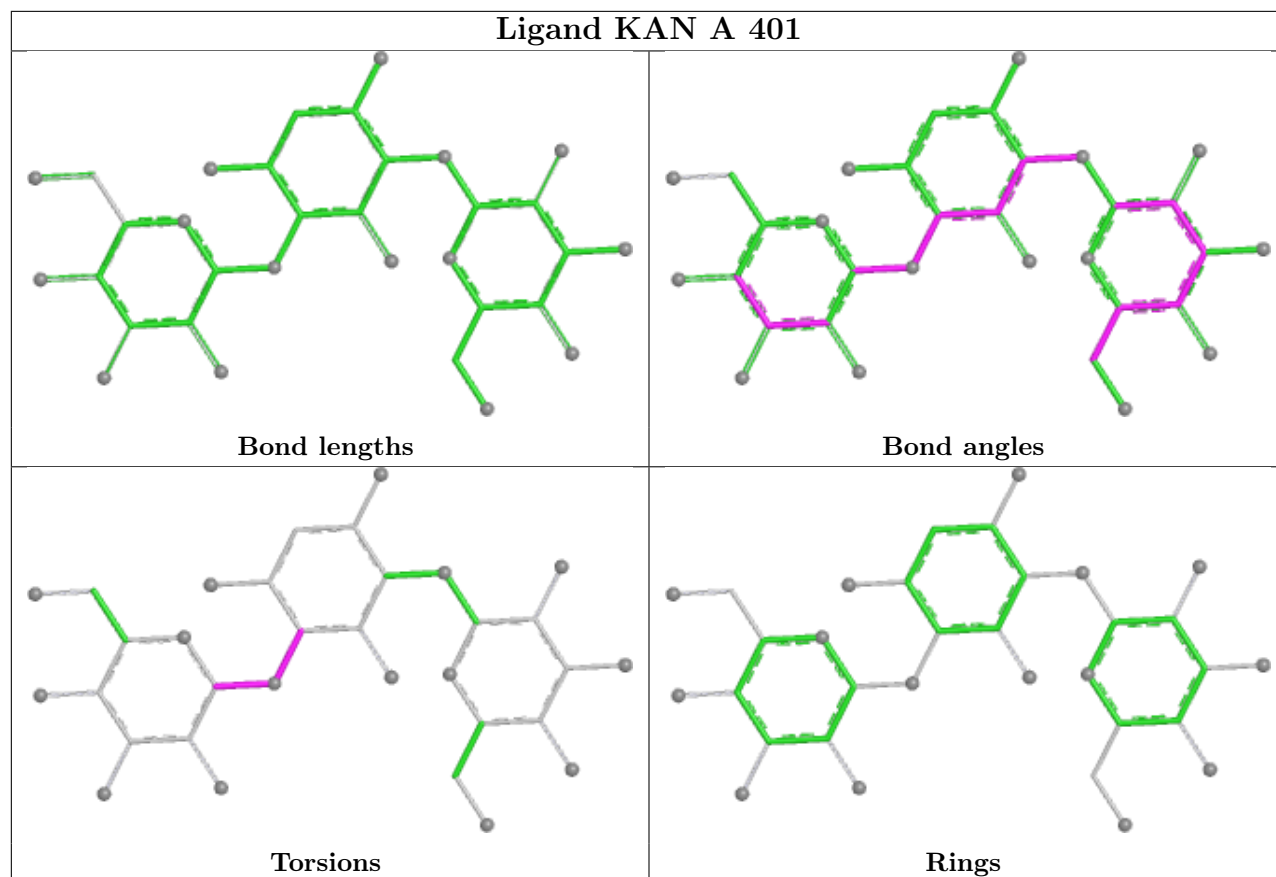
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	401[B]	QUE	5	0
3	A	402	QUE	2	0
3	B	401[A]	QUE	2	0
2	A	401	KAN	1	0

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









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	289/322 (89%)	0.80	23 (7%) 18 25	30, 66, 87, 113	3 (1%)
1	B	290/322 (90%)	0.78	19 (6%) 24 32	50, 65, 101, 118	0
All	All	579/644 (89%)	0.79	42 (7%) 21 28	30, 65, 94, 118	3 (0%)

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	140	SER	4.4
1	B	278	TYR	4.4
1	A	298	ILE	4.0
1	B	288	TYR	3.7
1	A	143	VAL	3.2
1	B	287	TRP	3.0
1	B	135	ILE	3.0
1	B	185	ILE	3.0
1	A	140	SER	2.9
1	B	298	ILE	2.9
1	B	282	TYR	2.9
1	B	279	GLY	2.6
1	B	142	LEU	2.6
1	B	295	ILE	2.6
1	A	99	LYS	2.6
1	A	190	TYR	2.6
1	A	287	TRP	2.6
1	A	236	ASP	2.6
1	A	141	ASN	2.4
1	A	284	TYR	2.4
1	A	3	THR	2.4
1	B	34	CYS	2.4
1	B	284	TYR	2.4
1	A	138	PHE	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	142	LEU	2.3
1	B	283	GLY	2.2
1	A	201	ASP	2.2
1	A	135	ILE	2.2
1	A	260	VAL	2.2
1	B	281	GLU	2.1
1	A	248	ILE	2.1
1	A	245	VAL	2.1
1	A	235	GLU	2.1
1	A	69	ASN	2.1
1	A	278	TYR	2.1
1	B	190	TYR	2.1
1	B	106	LEU	2.1
1	A	187	PHE	2.1
1	B	207	THR	2.1
1	A	257	ILE	2.0
1	A	113	LYS	2.0
1	B	134	ASN	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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

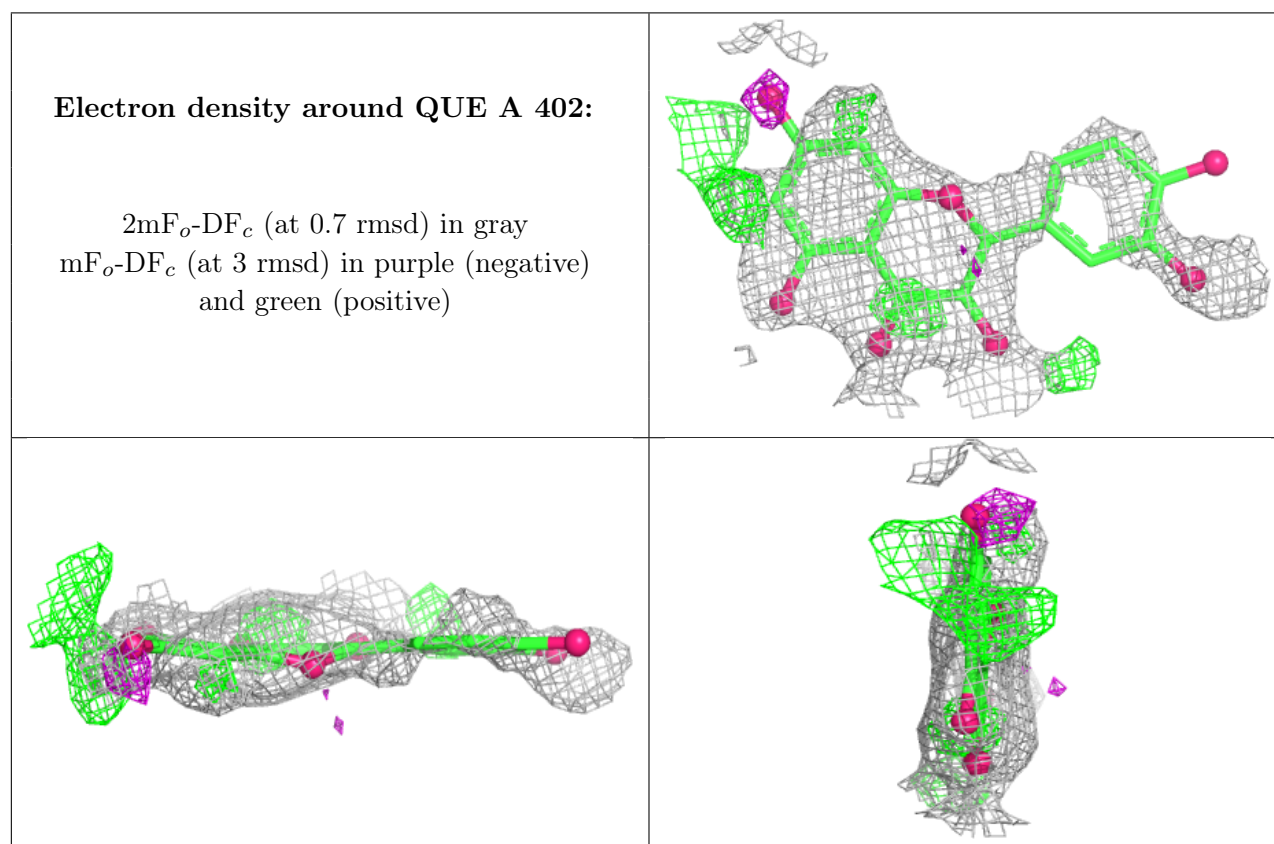
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	QUE	A	402	22/22	0.70	0.20	71,72,72,73	22
3	QUE	B	403	22/22	0.79	0.19	81,82,83,83	22
2	KAN	A	401	33/33	0.88	0.12	69,71,75,75	0
3	QUE	B	401[A]	22/22	0.88	0.15	70,71,71,71	22
3	QUE	B	401[B]	22/22	0.88	0.15	70,71,71,71	22

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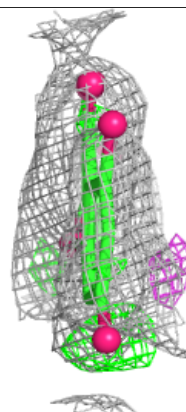
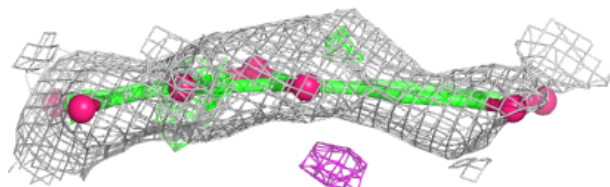
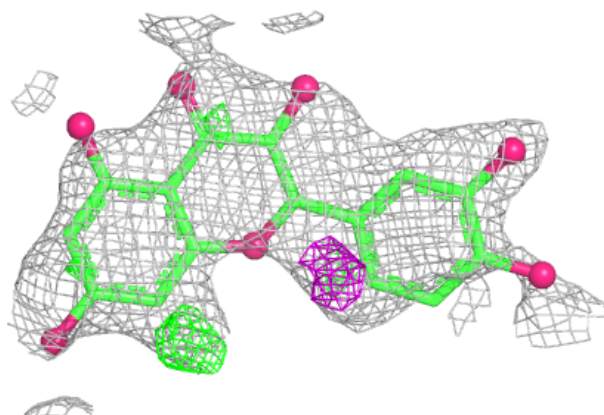
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	KAN	B	402	33/33	0.88	0.11	67,68,71,71	0
4	CL	B	404	1/1	0.91	0.11	86,86,86,86	0
4	CL	A	403	1/1	0.97	0.08	64,64,64,64	0

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.

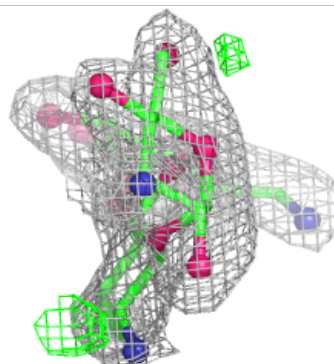
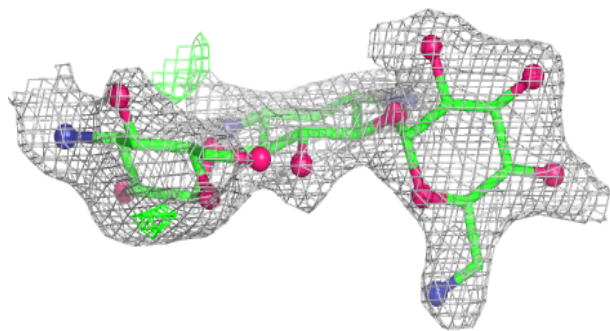
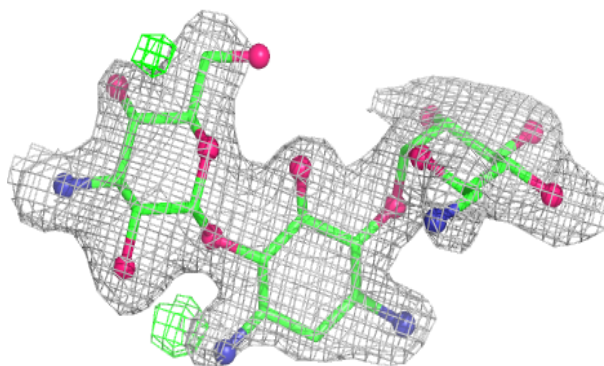


Electron density around QUE B 403:

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

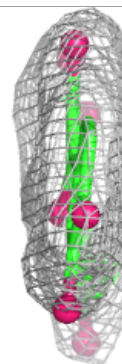
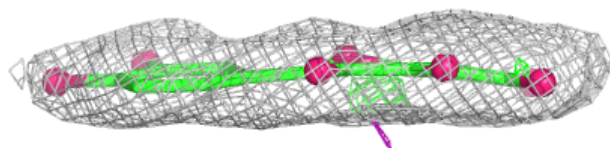
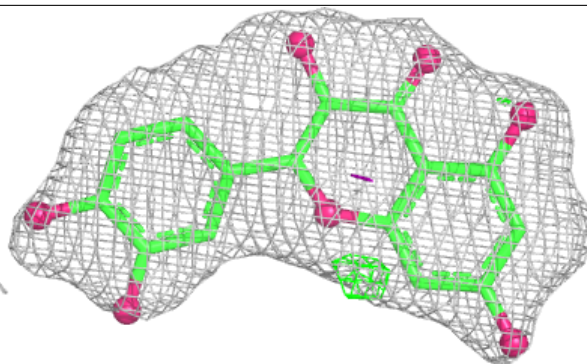
**Electron density around KAN A 401:**

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

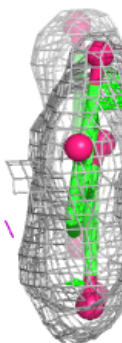
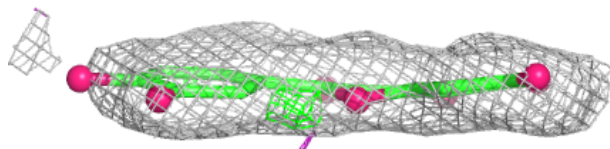
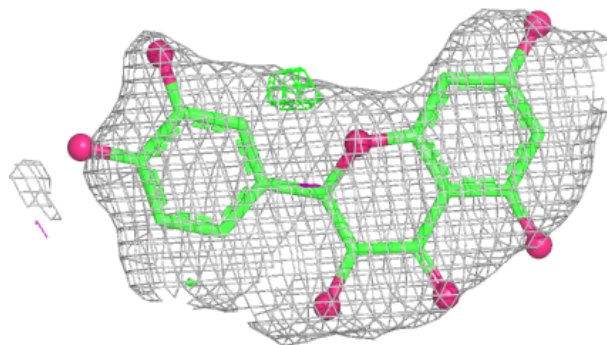


Electron density around QUE B 401 (A):

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

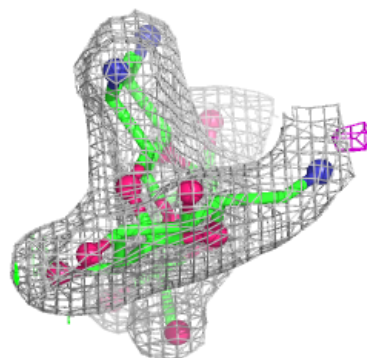
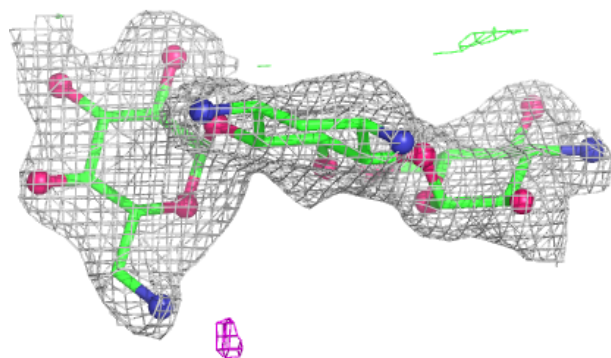
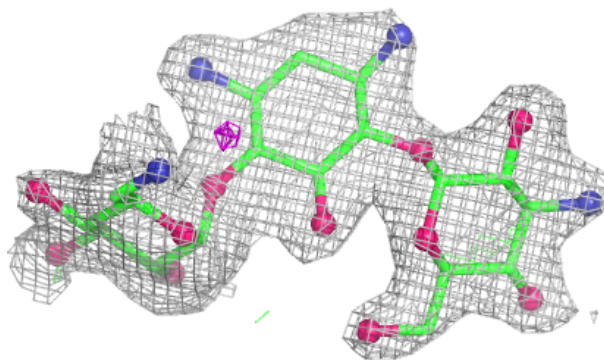
**Electron density around QUE B 401 (B):**

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



Electron density around KAN B 402:

$2mF_o-DF_c$ (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.