



# wwPDB X-ray Structure Validation Summary Report ⓘ

May 25, 2020 – 10:26 am BST

PDB ID : 6U7A  
Title : Rv3722c in complex with kynurenine  
Authors : Mandyoli, L.; Sacchettini, J.  
Deposited on : 2019-09-01  
Resolution : 2.22 Å(reported)

This is a wwPDB X-ray Structure Validation Summary Report for a publicly released PDB entry.

We welcome your comments at [validation@mail.wwpdb.org](mailto: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.

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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

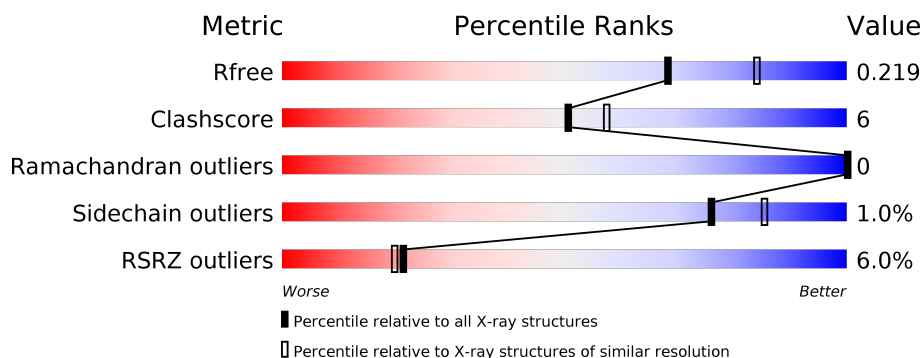
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.22 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	5912 (2.24-2.20)
Clashscore	141614	6646 (2.24-2.20)
Ramachandran outliers	138981	6543 (2.24-2.20)
Sidechain outliers	138945	6544 (2.24-2.20)
RSRZ outliers	127900	5797 (2.24-2.20)

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  $\geq 3$ , 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	435	<div> <div>4%</div> <div> <div></div> <div>89%</div> <div>8%</div> <div></div> </div> <div></div> </div>
1	B	435	<div> <div>5%</div> <div> <div></div> <div>85%</div> <div>12%</div> <div></div> </div> <div></div> </div>
1	C	435	<div> <div>7%</div> <div> <div></div> <div>87%</div> <div>10%</div> <div></div> </div> <div></div> </div>
1	D	435	<div> <div>4%</div> <div> <div></div> <div>88%</div> <div>9%</div> <div></div> </div> <div></div> </div>
1	E	435	<div> <div>7%</div> <div> <div></div> <div>88%</div> <div>9%</div> <div></div> </div> <div></div> </div>
1	F	435	<div> <div>2%</div> <div> <div></div> <div>87%</div> <div>10%</div> <div></div> </div> <div></div> </div>

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Mol	Chain	Length	Quality of chain
1	G	435	
1	H	435	

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
4	GOL	B	503	-	-	X	-
6	KYA	B	501	-	-	X	-
6	KYA	C	501	-	-	X	-
6	KYA	D	501	-	-	X	-
6	KYA	G	501	-	-	X	-
7	PMP	B	502	-	-	X	-

## 2 Entry composition

There are 9 unique types of molecules in this entry. The entry contains 27421 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.

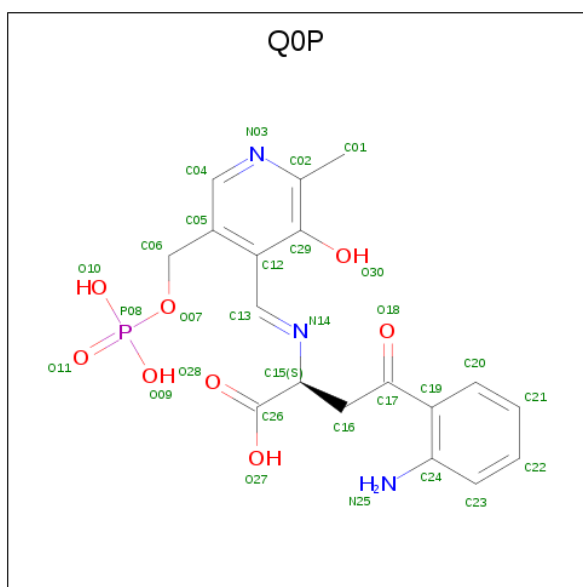
- Molecule 1 is a protein called Aminotransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	423	Total	C	N	O	S	0	0	0
			3255	2078	562	604	11			
1	B	423	Total	C	N	O	S	0	0	0
			3255	2078	562	604	11			
1	C	423	Total	C	N	O	S	0	0	0
			3255	2078	562	604	11			
1	D	429	Total	C	N	O	S	0	0	0
			3293	2100	569	613	11			
1	E	423	Total	C	N	O	S	0	0	0
			3255	2078	562	604	11			
1	F	422	Total	C	N	O	S	0	0	0
			3249	2075	561	602	11			
1	G	423	Total	C	N	O	S	0	0	0
			3255	2078	562	604	11			
1	H	416	Total	C	N	O	S	0	0	0
			3202	2046	554	591	11			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	VAL	-	expression tag	UNP A0A0E8TWE4
B	1	VAL	-	expression tag	UNP A0A0E8TWE4
C	1	VAL	-	expression tag	UNP A0A0E8TWE4
D	1	VAL	-	expression tag	UNP A0A0E8TWE4
E	1	VAL	-	expression tag	UNP A0A0E8TWE4
F	1	VAL	-	expression tag	UNP A0A0E8TWE4
G	1	VAL	-	expression tag	UNP A0A0E8TWE4
H	1	VAL	-	expression tag	UNP A0A0E8TWE4

- Molecule 2 is (2S)-4-(2-aminophenyl)-2-[(E)-({3-hydroxy-2-methyl-5-[(phosphonooxy)methyl]pyridin-4-yl}methylidene)amino]-4-oxobutanoic acid (three-letter code: Q0P) (formula: C<sub>18</sub>H<sub>20</sub>N<sub>3</sub>O<sub>8</sub>P) (labeled as "Ligand of Interest" by author).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			30	18	3	8	1		

- Molecule 3 is PHOSPHATE ION (three-letter code: PO4) (formula: O<sub>4</sub>P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	P	0	0
			5	4	1		
3	C	1	Total	O	P	0	0
			5	4	1		
3	D	1	Total	O	P	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	G	1	Total	O	P	0	0
			5	4	1		
3	G	1	Total	O	P	0	0
			5	4	1		
3	H	1	Total	O	P	0	0
			5	4	1		

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula:  $C_3H_8O_3$ ).



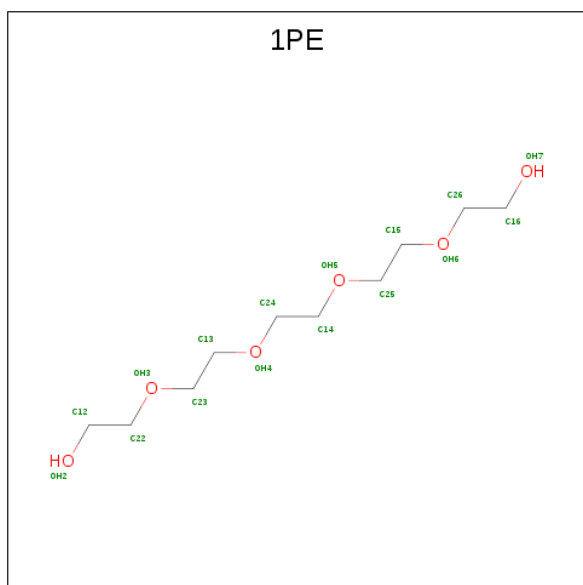
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			6	3	3		
4	A	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		
4	C	1	Total	C	O	0	0
			6	3	3		
4	D	1	Total	C	O	0	0
			6	3	3		
4	G	1	Total	C	O	0	0
			6	3	3		
4	G	1	Total	C	O	0	0
			6	3	3		

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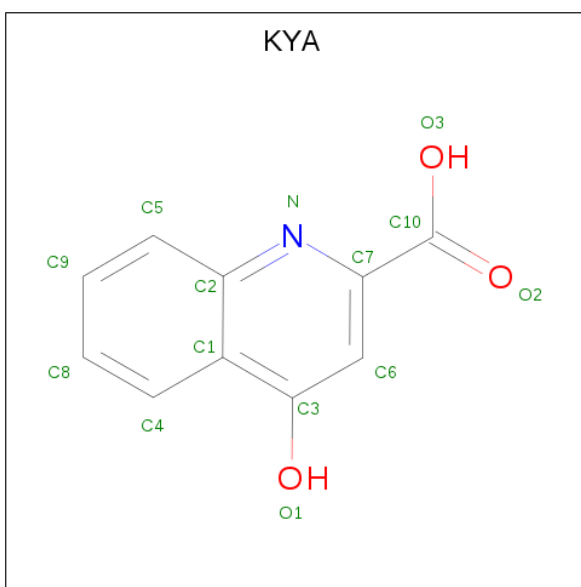
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	G	1	Total	C	O	0	0
			6	3	3		
4	G	1	Total	C	O	0	0
			6	3	3		

- Molecule 5 is PENTAETHYLENE GLYCOL (three-letter code: 1PE) (formula: C<sub>10</sub>H<sub>22</sub>O<sub>6</sub>).



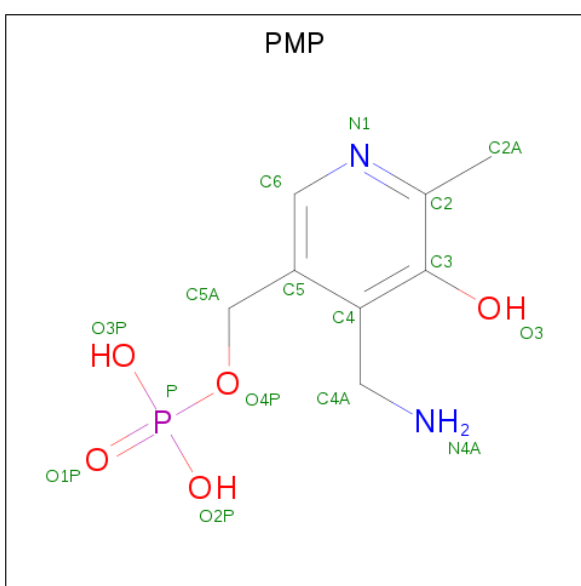
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			10	6	4		
5	B	1	Total	C	O	0	0
			16	10	6		
5	D	1	Total	C	O	0	0
			16	10	6		
5	E	1	Total	C	O	0	0
			10	6	4		
5	E	1	Total	C	O	0	0
			10	6	4		
5	H	1	Total	C	O	0	0
			7	4	3		

- Molecule 6 is 4-hydroxyquinoline-2-carboxylic acid (three-letter code: KYA) (formula: C<sub>10</sub>H<sub>7</sub>NO<sub>3</sub>) (labeled as "Ligand of Interest" by author).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	B	1	Total	C	N	O	0	0
			14	10	1	3		
6	C	1	Total	C	N	O	0	0
			14	10	1	3		
6	D	1	Total	C	N	O	0	0
			14	10	1	3		
6	G	1	Total	C	N	O	0	0
			14	10	1	3		

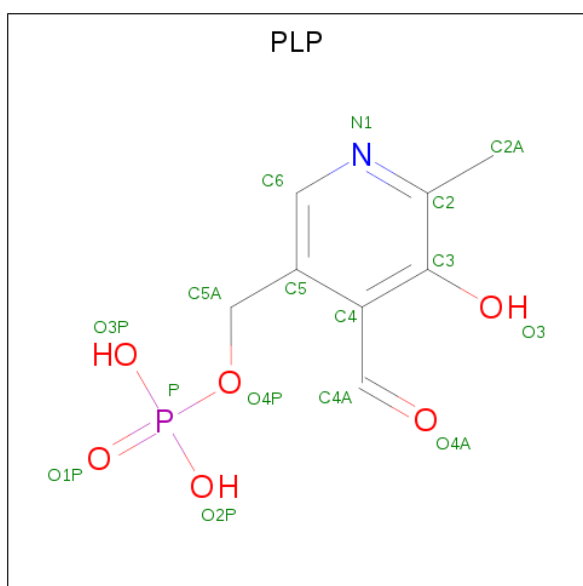
- Molecule 7 is 4'-DEOXY-4'-AMINOPYRIDOXAL-5'-PHOSPHATE (three-letter code: PMP) (formula:  $C_8H_{13}N_2O_5P$ ).





Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
7	B	1	Total	C	N	O	P	0	0
			16	8	2	5	1		
7	C	1	Total	C	N	O	P	0	0
			16	8	2	5	1		
7	D	1	Total	C	N	O	P	0	0
			16	8	2	5	1		
7	G	1	Total	C	N	O	P	0	0
			16	8	2	5	1		

- Molecule 8 is PYRIDOXAL-5'-PHOSPHATE (three-letter code: PLP) (formula:  $C_8H_{10}NO_6P$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
8	E	1	Total	C	N	O	P	0	0
			16	8	1	6	1		
8	F	1	Total	C	N	O	P	0	0
			16	8	1	6	1		
8	H	1	Total	C	N	O	P	0	0
			16	8	1	6	1		

- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	145	Total	O	0	0
			145	145		
9	B	122	Total	O	0	0
			122	122		
9	C	136	Total	O	0	0
			136	136		

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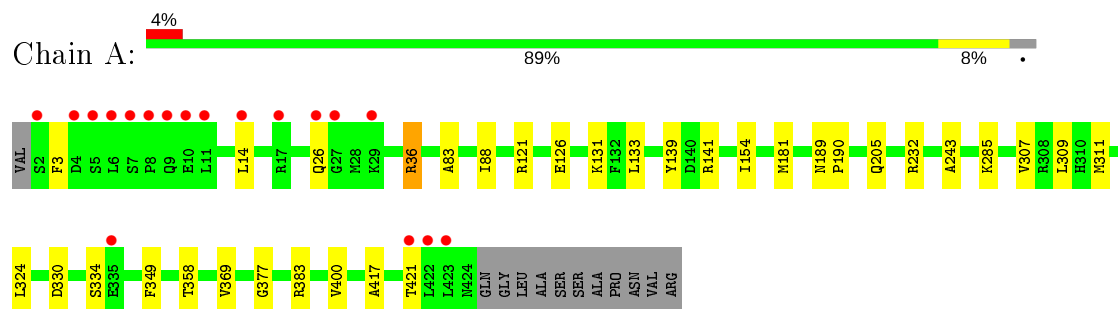
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	D	144	Total 144	O 144	0	0
9	E	118	Total 118	O 118	0	0
9	F	145	Total 145	O 145	0	0
9	G	122	Total 122	O 122	0	0
9	H	113	Total 113	O 113	0	0

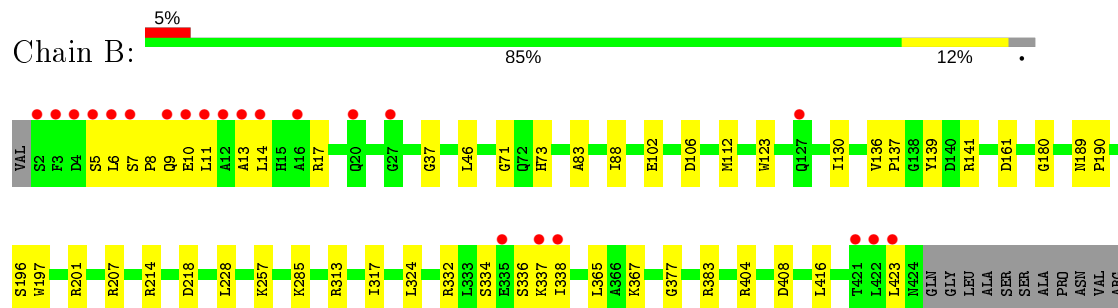
### 3 Residue-property plots [i](#)

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.

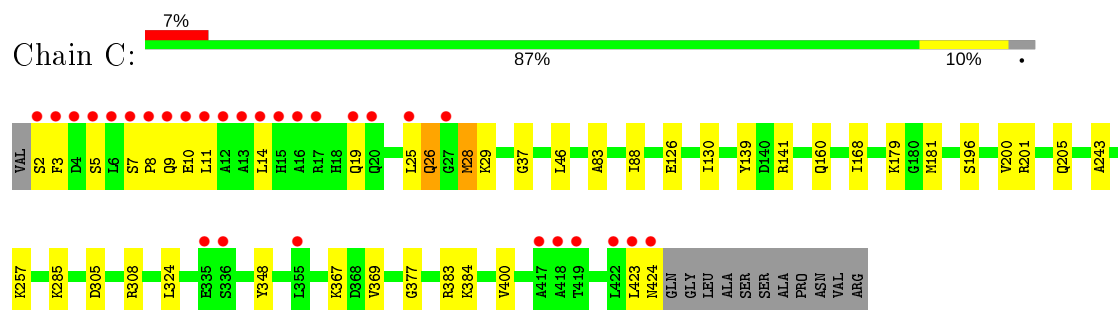
- Molecule 1: Aminotransferase



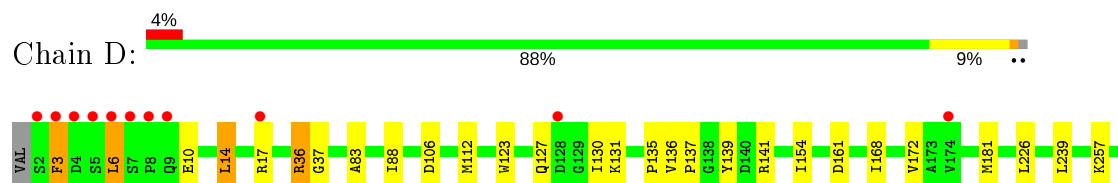
- Molecule 1: Aminotransferase

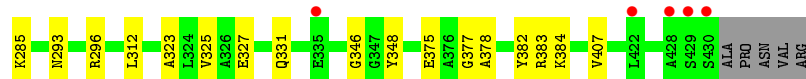


- Molecule 1: Aminotransferase

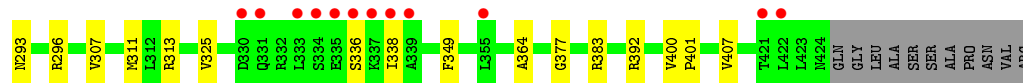
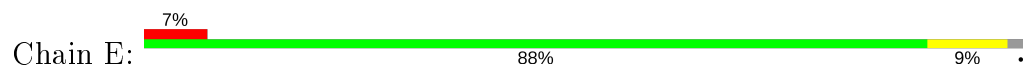


- Molecule 1: Aminotransferase

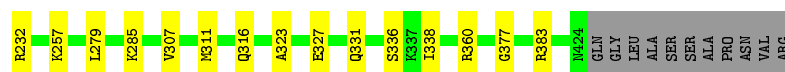
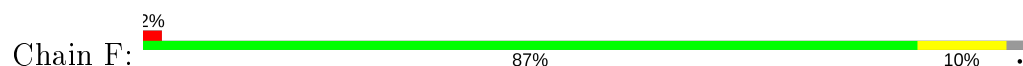




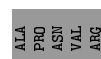
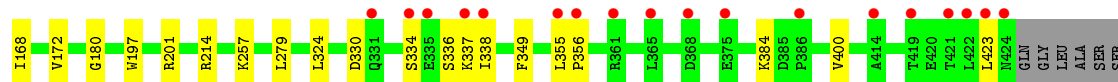
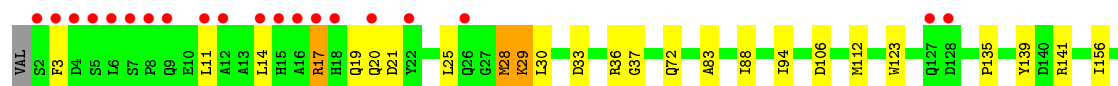
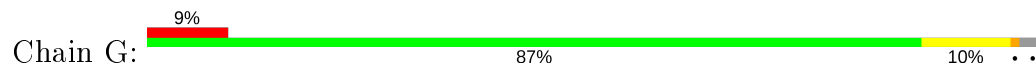
- Molecule 1: Aminotransferase



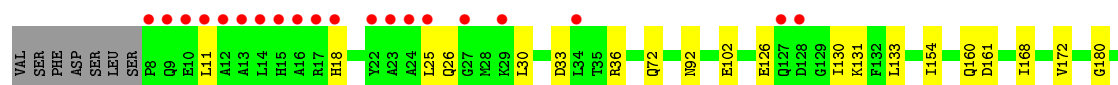
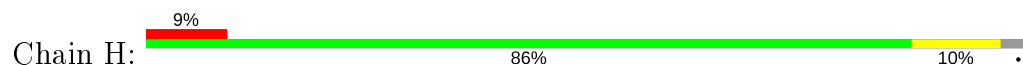
- Molecule 1: Aminotransferase









- Molecule 1: Aminotransferase



- Molecule 1: Aminotransferase



A417	A418	T419	E420	T421	L422	L423	ASN	GLN	GLY	LEU	ALA	SER	SER	ALA	PRO	ASN	VAL	ARG
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## 4 Data and refinement statistics

Property	Value	Source
Space group	P 31	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	108.40Å 108.40Å 321.34Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	32.46 – 2.22 48.36 – 2.22	Depositor EDS
% Data completeness (in resolution range)	85.2 (32.46-2.22) 85.3 (48.36-2.22)	Depositor EDS
$R_{merge}$	0.14	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.74 (at 2.22Å)	Xtriage
Refinement program	PHENIX 1.15.2_3472, PHENIX 1.15.2_3472	Depositor
R, $R_{free}$	0.178 , 0.219 0.179 , 0.219	Depositor DCC
$R_{free}$ test set	8912 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	31.1	Xtriage
Anisotropy	0.007	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.38 , 48.4	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.003 for -h,-k,l 0.027 for h,-h-k,-l 0.017 for -k,-h,-l	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	27421	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.35% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: GOL, PMP, PLP, PO4, 1PE, Q0P, KYA

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	0.52	0/3335	0.74	0/4540
1	B	0.50	0/3335	0.74	0/4540
1	C	0.51	0/3335	0.72	0/4540
1	D	0.47	0/3373	0.73	0/4591
1	E	0.46	0/3335	0.74	0/4540
1	F	0.46	0/3329	0.70	0/4532
1	G	0.46	0/3335	0.73	0/4540
1	H	0.46	0/3281	0.73	0/4466
All	All	0.48	0/26658	0.73	0/36289

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3255	0	3228	24	0
1	B	3255	0	3228	50	0
1	C	3255	0	3228	37	0
1	D	3293	0	3265	38	0
1	E	3255	0	3228	46	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	F	3249	0	3223	26	0
1	G	3255	0	3228	45	0
1	H	3202	0	3184	26	0
2	A	30	0	0	2	0
3	A	5	0	0	0	0
3	C	5	0	0	0	0
3	D	5	0	0	0	0
3	G	10	0	0	0	0
3	H	5	0	0	0	0
4	A	12	0	16	1	0
4	B	12	0	16	5	0
4	C	6	0	8	2	0
4	D	6	0	8	2	0
4	G	24	0	31	0	0
5	A	10	0	12	2	0
5	B	16	0	22	5	0
5	D	16	0	22	4	0
5	E	20	0	24	4	0
5	H	7	0	8	2	0
6	B	14	0	6	16	0
6	C	14	0	5	10	0
6	D	14	0	5	11	0
6	G	14	0	6	15	0
7	B	16	0	10	6	0
7	C	16	0	11	3	0
7	D	16	0	11	3	0
7	G	16	0	11	3	0
8	E	16	0	8	1	0
8	F	16	0	8	1	0
8	H	16	0	8	0	0
9	A	145	0	0	2	0
9	B	122	0	0	2	0
9	C	136	0	0	0	0
9	D	144	0	0	0	0
9	E	118	0	0	0	0
9	F	145	0	0	2	0
9	G	122	0	0	1	0
9	H	113	0	0	0	0
All	All	27421	0	26068	300	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 6.



The worst 5 of 300 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:C:37:GLY:O	6:C:501:KYA:H5	1.61	0.97
1:D:141:ARG:HH21	6:D:501:KYA:H6	1.25	0.97
1:B:37:GLY:O	6:B:501:KYA:H5	1.65	0.96
1:B:257:LYS:NZ	6:B:501:KYA:H9	1.83	0.92
1:G:28:MET:HE2	1:G:30:LEU:HD21	1.57	0.85

There are no symmetry-related clashes.

## 5.3 Torsion angles [i](#)

### 5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	421/435 (97%)	411 (98%)	10 (2%)	0	100	100
1	B	421/435 (97%)	412 (98%)	9 (2%)	0	100	100
1	C	421/435 (97%)	414 (98%)	7 (2%)	0	100	100
1	D	427/435 (98%)	419 (98%)	8 (2%)	0	100	100
1	E	421/435 (97%)	413 (98%)	8 (2%)	0	100	100
1	F	420/435 (97%)	411 (98%)	9 (2%)	0	100	100
1	G	421/435 (97%)	412 (98%)	9 (2%)	0	100	100
1	H	414/435 (95%)	406 (98%)	8 (2%)	0	100	100
All	All	3366/3480 (97%)	3298 (98%)	68 (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	343/352 (97%)	342 (100%)	1 (0%)	92	96
1	B	343/352 (97%)	342 (100%)	1 (0%)	92	96
1	C	343/352 (97%)	337 (98%)	6 (2%)	60	73
1	D	347/352 (99%)	342 (99%)	5 (1%)	67	78
1	E	343/352 (97%)	340 (99%)	3 (1%)	78	87
1	F	342/352 (97%)	340 (99%)	2 (1%)	86	92
1	G	343/352 (97%)	339 (99%)	4 (1%)	71	82
1	H	336/352 (96%)	331 (98%)	5 (2%)	65	76
All	All	2740/2816 (97%)	2713 (99%)	27 (1%)	76	85

5 of 27 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	36	ARG
1	E	207	ARG
1	H	207	ARG
1	E	6	LEU
1	C	28	MET

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	E	205	GLN
1	H	19	GLN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

34 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
4	GOL	A	504	-	5,5,5	0.81	0	5,5,5	0.95	0
3	PO4	G	503	-	4,4,4	1.03	0	6,6,6	0.56	0
6	KYA	B	501	-	12,15,15	2.63	6 (50%)	16,21,21	2.88	6 (37%)
3	PO4	H	502	-	4,4,4	0.85	0	6,6,6	0.53	0
6	KYA	G	501	-	12,15,15	1.98	4 (33%)	16,21,21	1.55	5 (31%)
5	1PE	D	505	-	15,15,15	0.53	0	14,14,14	0.39	0
4	GOL	B	504	-	5,5,5	1.36	0	5,5,5	0.80	0
4	GOL	A	503	-	5,5,5	1.09	0	5,5,5	0.75	0
6	KYA	C	501	-	12,15,15	2.59	3 (25%)	16,21,21	2.49	3 (18%)
4	GOL	C	504	-	5,5,5	1.08	0	5,5,5	0.63	0
3	PO4	C	503	-	4,4,4	0.96	0	6,6,6	0.70	0
3	PO4	A	502	-	4,4,4	1.16	0	6,6,6	0.82	0
4	GOL	B	503	-	5,5,5	1.18	0	5,5,5	1.01	0
7	PMP	G	502	-	16,16,16	0.95	1 (6%)	21,23,23	1.34	2 (9%)
5	1PE	H	503	-	6,6,15	0.47	0	5,5,14	0.38	0
8	PLP	F	501	-	16,16,16	1.20	2 (12%)	20,23,23	1.34	4 (20%)
3	PO4	D	503	-	4,4,4	0.85	0	6,6,6	0.91	0
5	1PE	B	505	-	15,15,15	0.52	0	14,14,14	0.63	0
4	GOL	G	507	-	5,5,5	1.45	1 (20%)	5,5,5	0.83	0
7	PMP	B	502	-	16,16,16	2.88	3 (18%)	21,23,23	1.21	1 (4%)
6	KYA	D	501	-	12,15,15	3.11	6 (50%)	16,21,21	2.96	4 (25%)
5	1PE	A	505	-	9,9,15	0.47	0	8,8,14	0.68	0
7	PMP	D	502	-	16,16,16	0.95	1 (6%)	21,23,23	1.45	2 (9%)
5	1PE	E	502	-	9,9,15	0.47	0	8,8,14	0.77	0
2	Q0P	A	501	-	27,31,31	1.69	5 (18%)	34,44,44	2.14	9 (26%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
4	GOL	G	505	-	5,5,5	1.38	1 (20%)	5,5,5	0.80	0
4	GOL	G	508	-	5,5,5	0.95	0	5,5,5	0.99	0
5	1PE	E	503	-	9,9,15	0.53	0	8,8,14	0.57	0
4	GOL	G	506	-	5,5,5	1.27	1 (20%)	5,5,5	1.05	0
4	GOL	D	504	-	5,5,5	1.12	0	5,5,5	0.92	0
8	PLP	E	501	-	16,16,16	1.21	2 (12%)	20,23,23	1.69	4 (20%)
8	PLP	H	501	-	16,16,16	1.18	2 (12%)	20,23,23	1.30	2 (10%)
7	PMP	C	502	-	16,16,16	1.04	1 (6%)	21,23,23	1.38	2 (9%)
3	PO4	G	504	-	4,4,4	0.90	0	6,6,6	0.65	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	GOL	A	504	-	-	0/4/4/4	-
5	1PE	H	503	-	-	1/4/4/13	-
6	KYA	B	501	-	-	0/0/4/4	0/2/2/2
6	KYA	G	501	-	-	0/0/4/4	0/2/2/2
5	1PE	D	505	-	-	10/13/13/13	-
4	GOL	B	504	-	-	0/4/4/4	-
4	GOL	A	503	-	-	4/4/4/4	-
6	KYA	C	501	-	-	0/0/4/4	0/2/2/2
4	GOL	C	504	-	-	2/4/4/4	-
5	1PE	E	502	-	-	6/7/7/13	-
8	PLP	F	501	-	-	2/8/8/8	0/1/1/1
7	PMP	G	502	-	-	5/8/8/8	0/1/1/1
4	GOL	B	503	-	-	4/4/4/4	-
5	1PE	B	505	-	-	7/13/13/13	-
4	GOL	G	507	-	-	4/4/4/4	-
7	PMP	B	502	-	-	5/8/8/8	0/1/1/1
6	KYA	D	501	-	-	0/0/4/4	0/2/2/2
5	1PE	A	505	-	-	6/7/7/13	-
7	PMP	D	502	-	-	5/8/8/8	0/1/1/1
2	Q0P	A	501	-	-	11/19/23/23	0/2/2/2
4	GOL	G	505	-	-	4/4/4/4	-
4	GOL	G	508	-	-	2/4/4/4	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	1PE	E	503	-	-	5/7/7/13	-
4	GOL	G	506	-	-	3/4/4/4	-
4	GOL	D	504	-	-	0/4/4/4	-
8	PLP	E	501	-	-	2/8/8/8	0/1/1/1
8	PLP	H	501	-	-	2/8/8/8	0/1/1/1
7	PMP	C	502	-	-	5/8/8/8	0/1/1/1

The worst 5 of 39 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	B	502	PMP	C3-C2	7.99	1.48	1.40
6	D	501	KYA	C1-C2	-7.31	1.30	1.42
6	C	501	KYA	C1-C2	-6.73	1.31	1.42
7	B	502	PMP	C5-C4	5.89	1.48	1.40
2	A	501	Q0P	C12-C13	5.62	1.57	1.46

The worst 5 of 44 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	D	501	KYA	C7-N-C2	9.26	124.95	118.06
2	A	501	Q0P	C15-N14-C13	8.26	128.60	117.40
6	C	501	KYA	C7-N-C2	7.64	123.74	118.06
6	B	501	KYA	C7-N-C2	6.46	122.87	118.06
6	B	501	KYA	C6-C7-N	-5.70	117.03	122.23

There are no chirality outliers.

5 of 95 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	503	GOL	O1-C1-C2-C3
4	A	503	GOL	C1-C2-C3-O3
4	A	503	GOL	O2-C2-C3-O3
4	C	504	GOL	C1-C2-C3-O3
4	C	504	GOL	O2-C2-C3-O3

There are no ring outliers.

22 monomers are involved in 88 short contacts:

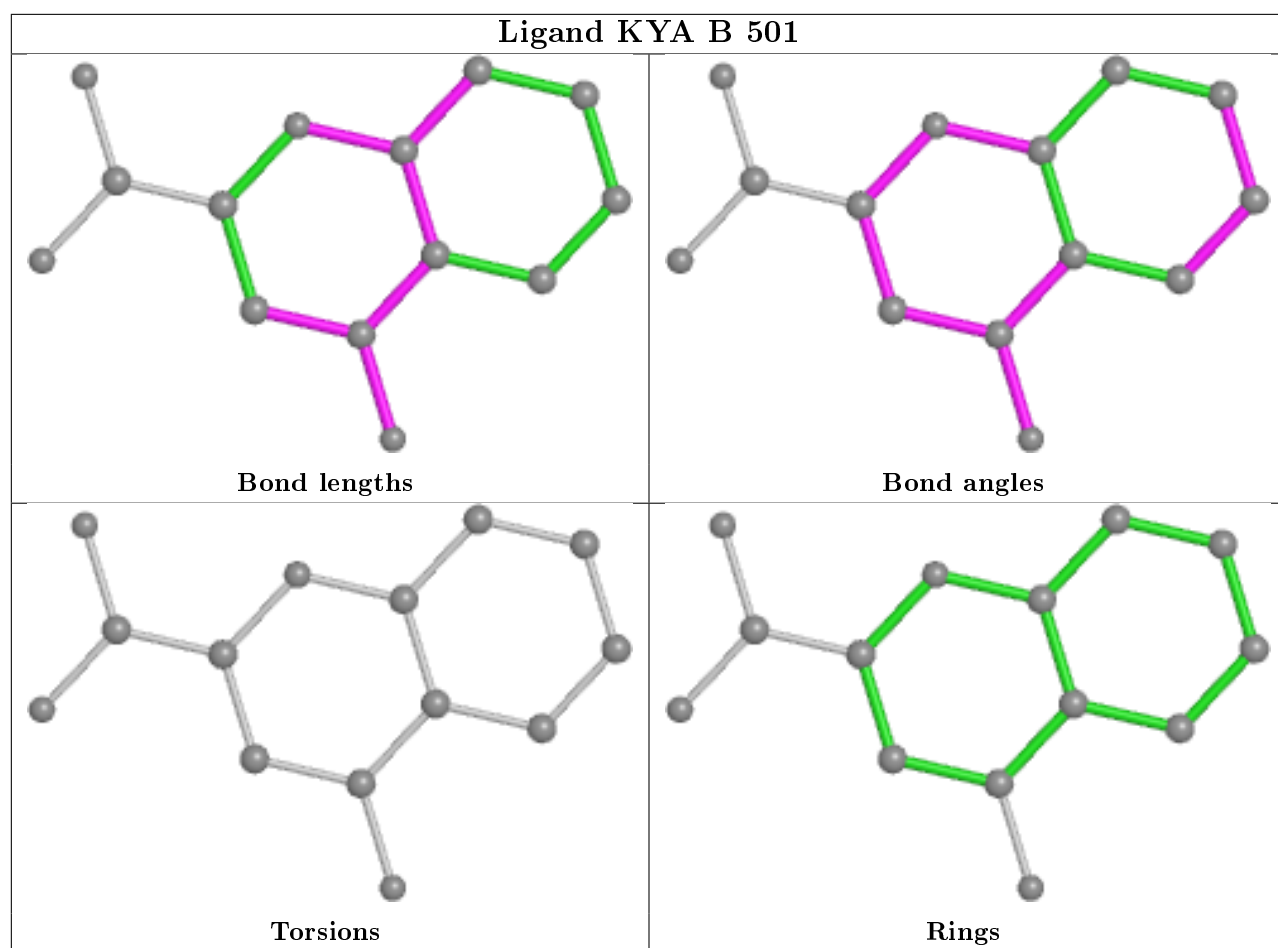
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	504	GOL	1	0

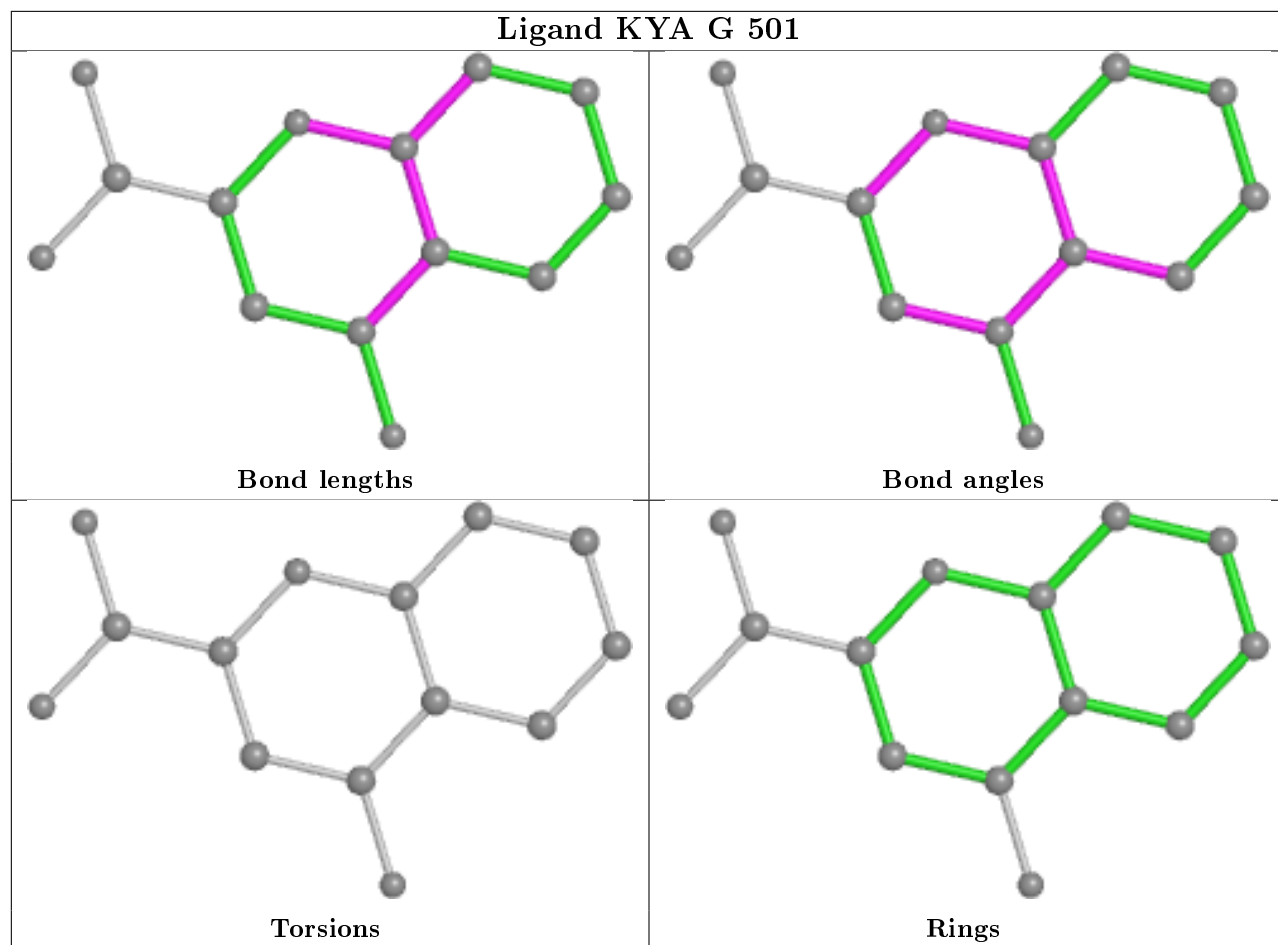
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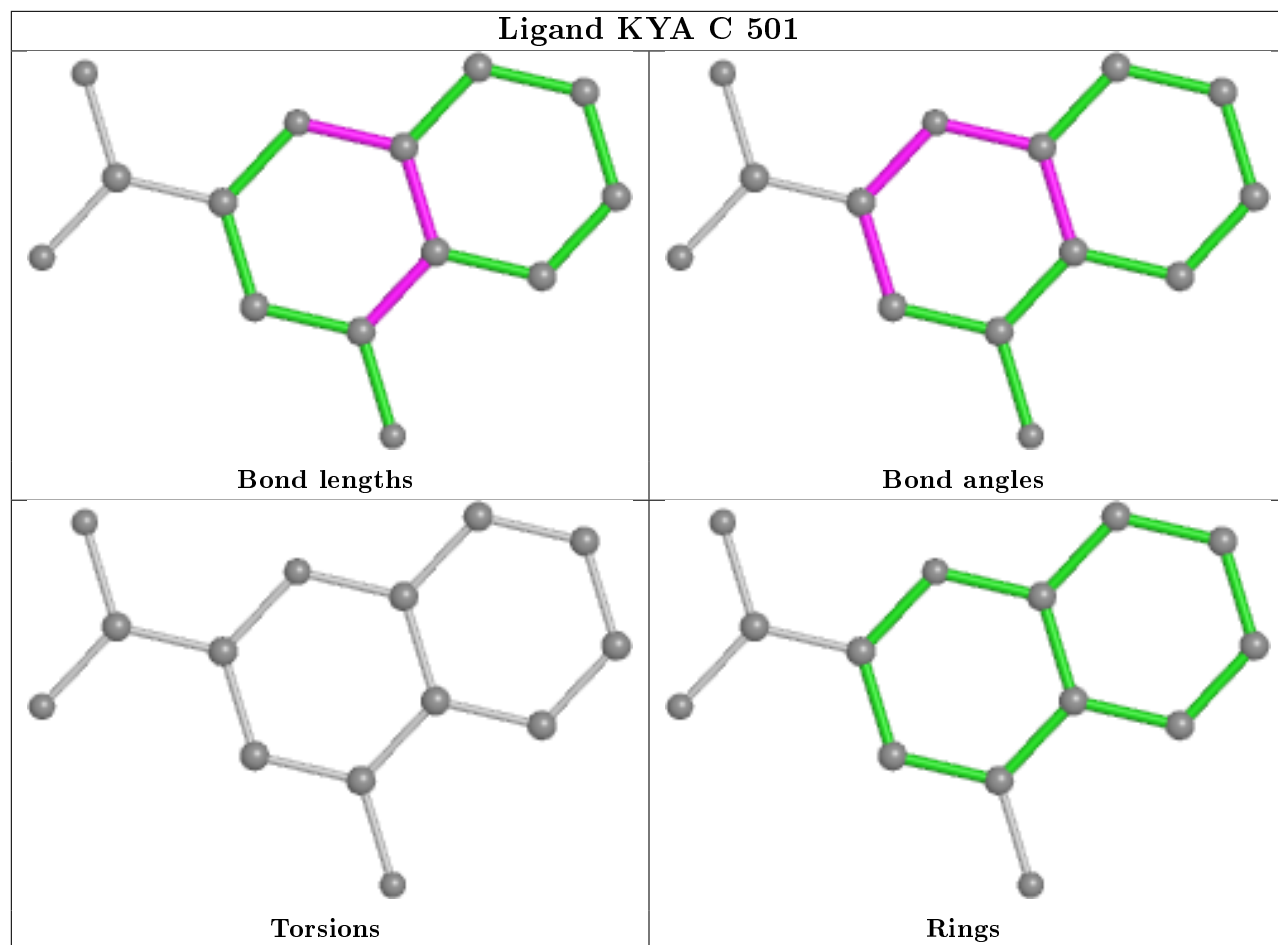
Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	B	501	KYA	16	0
6	G	501	KYA	15	0
5	D	505	1PE	4	0
4	B	504	GOL	1	0
6	C	501	KYA	10	0
4	C	504	GOL	2	0
4	B	503	GOL	4	0
7	G	502	PMP	3	0
5	H	503	1PE	2	0
8	F	501	PLP	1	0
5	B	505	1PE	5	0
7	B	502	PMP	6	0
6	D	501	KYA	11	0
5	A	505	1PE	2	0
7	D	502	PMP	3	0
5	E	502	1PE	1	0
2	A	501	Q0P	2	0
5	E	503	1PE	3	0
4	D	504	GOL	2	0
8	E	501	PLP	1	0
7	C	502	PMP	3	0

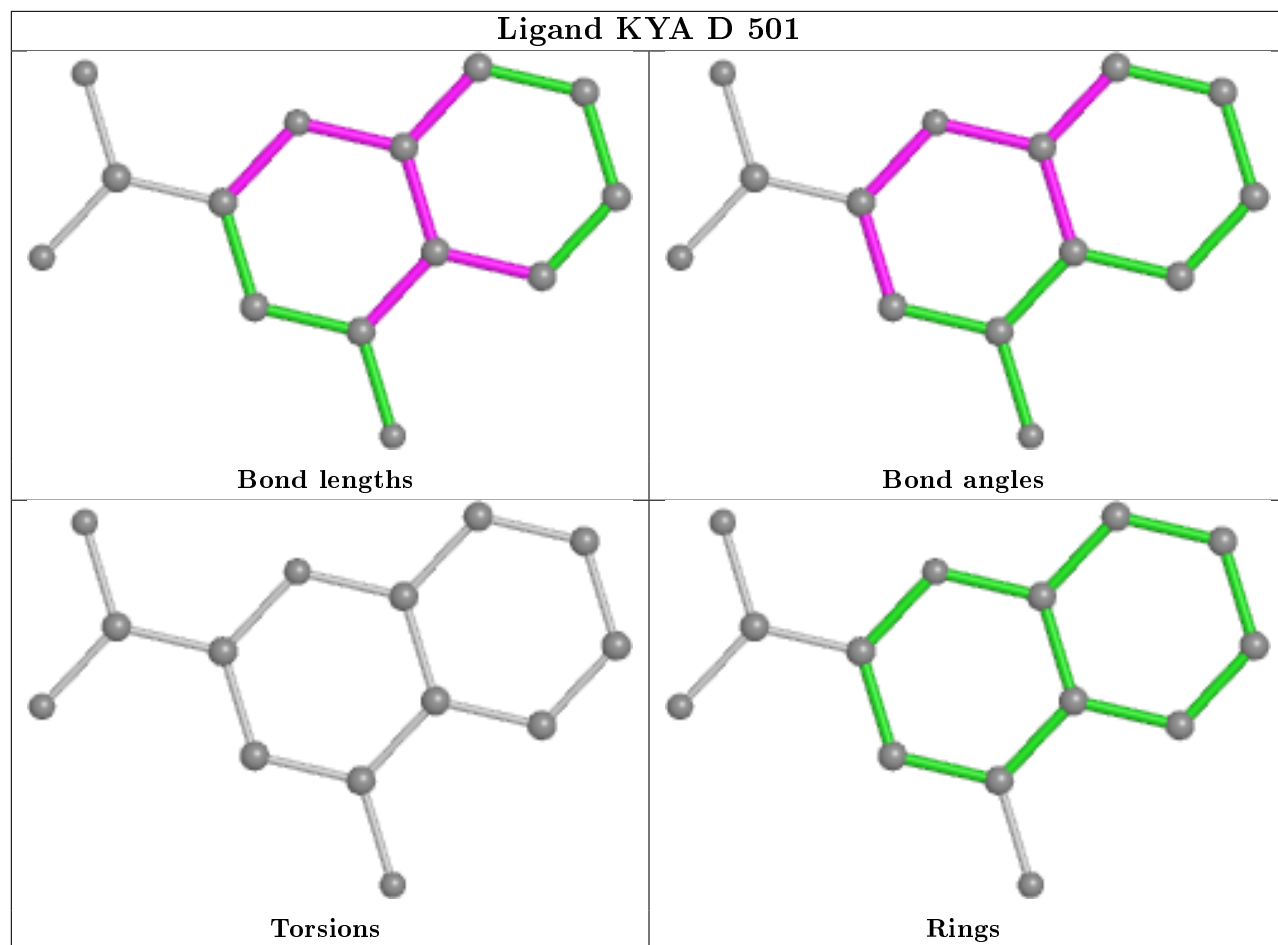
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

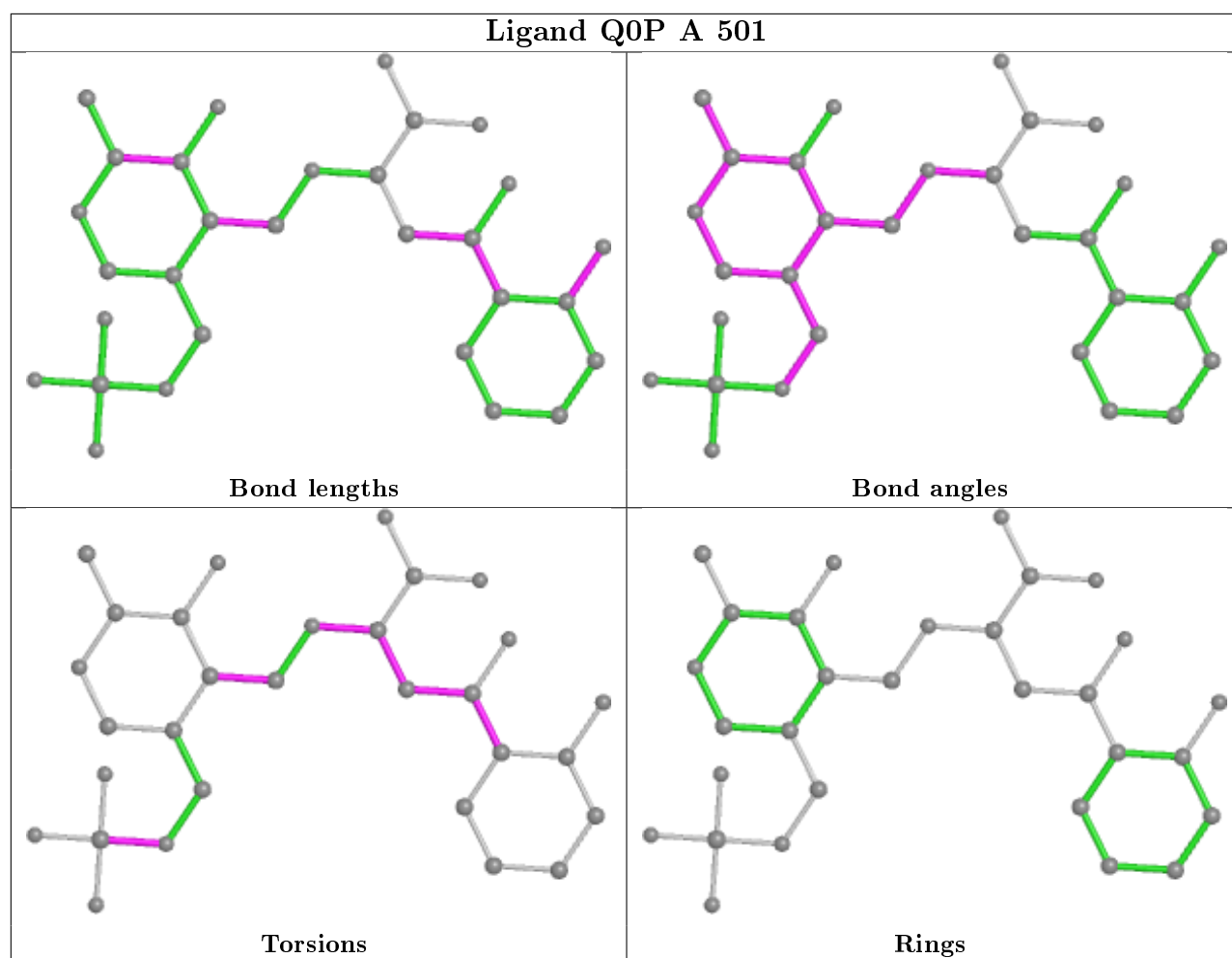












## 5.7 Other polymers [i](#)

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	423/435 (97%)	-0.21	18 (4%)	35	33	18, 28, 57, 78	1 (0%)
1	B	423/435 (97%)	-0.16	22 (5%)	27	25	20, 32, 64, 86	0
1	C	423/435 (97%)	-0.08	29 (6%)	16	15	20, 32, 63, 90	0
1	D	429/435 (98%)	-0.18	16 (3%)	41	39	20, 33, 53, 81	1 (0%)
1	E	423/435 (97%)	0.00	30 (7%)	16	14	21, 35, 65, 88	0
1	F	422/435 (97%)	-0.14	10 (2%)	59	57	23, 33, 54, 83	0
1	G	423/435 (97%)	0.22	38 (8%)	9	8	22, 35, 66, 81	0
1	H	416/435 (95%)	0.19	41 (9%)	7	6	22, 36, 74, 95	0
All	All	3382/3480 (97%)	-0.05	204 (6%)	21	20	18, 33, 64, 95	2 (0%)

The worst 5 of 204 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	H	12	ALA	8.8
1	C	6	LEU	8.0
1	H	11	LEU	7.5
1	E	5	SER	6.8
1	A	5	SER	6.7

### 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 ⓘ

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, 95<sup>th</sup> 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.

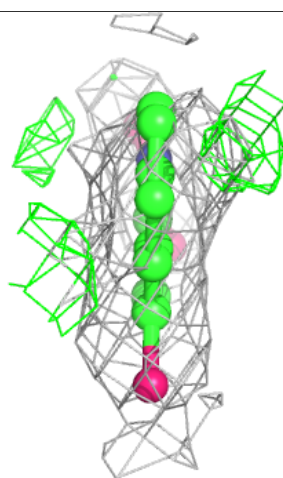
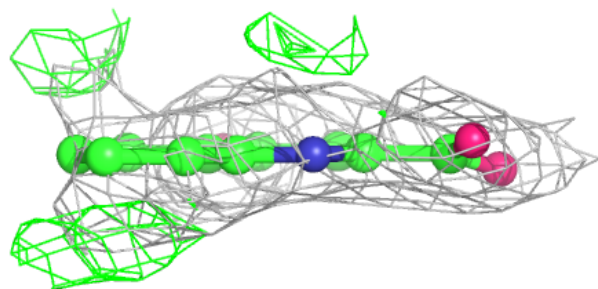
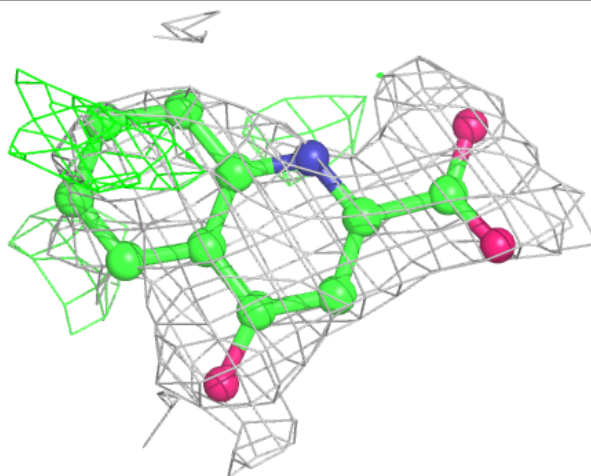
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
4	GOL	D	504	6/6	0.80	0.18	39,51,55,63	0
4	GOL	C	504	6/6	0.81	0.22	41,50,54,57	0
5	1PE	D	505	16/16	0.83	0.31	56,64,70,72	0
4	GOL	G	507	6/6	0.84	0.20	33,43,45,46	0
4	GOL	G	506	6/6	0.85	0.16	42,43,48,53	0
4	GOL	G	508	6/6	0.86	0.19	54,56,57,61	0
6	KYA	G	501	14/14	0.86	0.27	37,40,42,45	13
4	GOL	B	504	6/6	0.86	0.20	37,41,44,46	0
5	1PE	E	503	10/16	0.87	0.14	47,53,57,58	0
4	GOL	A	503	6/6	0.88	0.14	34,38,44,46	0
5	1PE	B	505	16/16	0.89	0.16	34,49,55,60	0
6	KYA	C	501	14/14	0.89	0.19	34,39,41,42	13
3	PO4	G	504	5/5	0.89	0.26	47,51,56,57	4
6	KYA	D	501	14/14	0.90	0.21	37,42,44,46	11
4	GOL	A	504	6/6	0.91	0.12	41,46,52,57	0
6	KYA	B	501	14/14	0.91	0.18	31,38,41,47	10
4	GOL	B	503	6/6	0.92	0.15	33,35,39,41	0
5	1PE	H	503	7/16	0.92	0.26	57,57,59,63	0
5	1PE	E	502	10/16	0.93	0.12	31,39,48,49	0
5	1PE	A	505	10/16	0.93	0.12	36,43,52,55	0
4	GOL	G	505	6/6	0.94	0.15	32,40,42,43	0
3	PO4	A	502	5/5	0.95	0.14	25,27,30,35	2
3	PO4	D	503	5/5	0.96	0.24	45,46,49,54	3
8	PLP	H	501	16/16	0.96	0.15	23,35,44,53	0
3	PO4	G	503	5/5	0.96	0.11	30,35,36,40	3
3	PO4	H	502	5/5	0.97	0.11	51,52,55,56	3
8	PLP	F	501	16/16	0.97	0.15	22,34,39,44	0
2	Q0P	A	501	30/30	0.97	0.16	20,29,39,43	17
8	PLP	E	501	16/16	0.97	0.12	26,37,44,52	0
7	PMP	B	502	16/16	0.97	0.15	20,32,36,39	0
7	PMP	C	502	16/16	0.97	0.14	24,32,38,38	0
7	PMP	G	502	16/16	0.97	0.14	22,37,41,48	0
7	PMP	D	502	16/16	0.98	0.12	23,33,40,43	0
3	PO4	C	503	5/5	0.98	0.13	38,38,42,43	4

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.

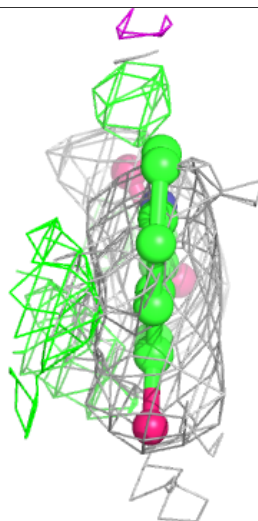
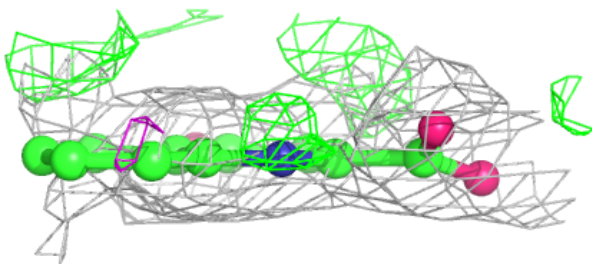
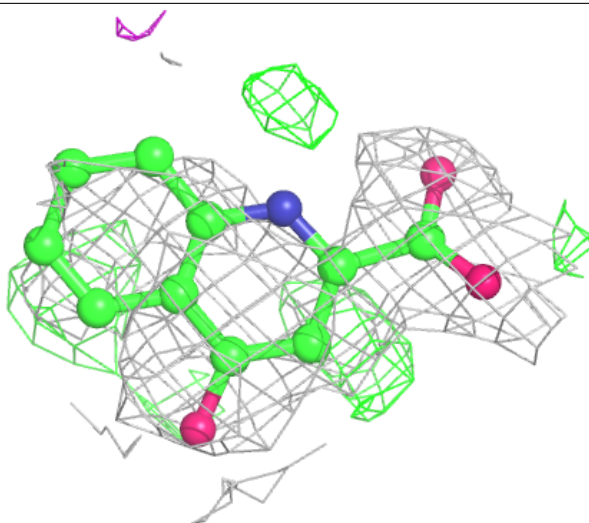
**Electron density around KYA G 501:**

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



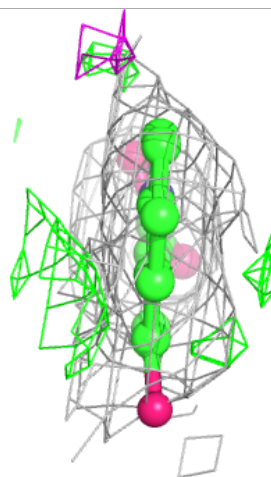
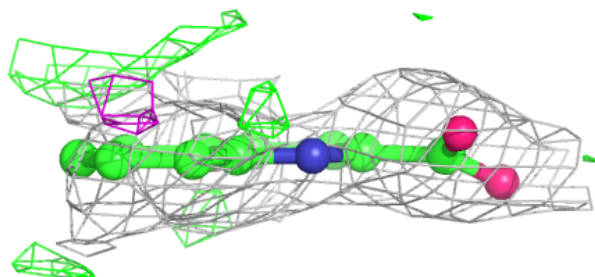
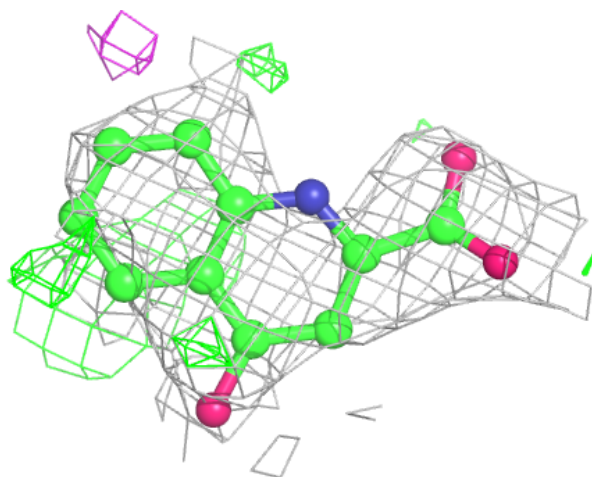
**Electron density around KYA C 501:**

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



**Electron density around KYA D 501:**

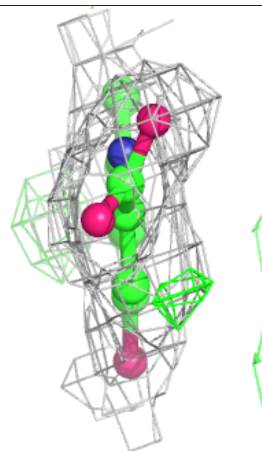
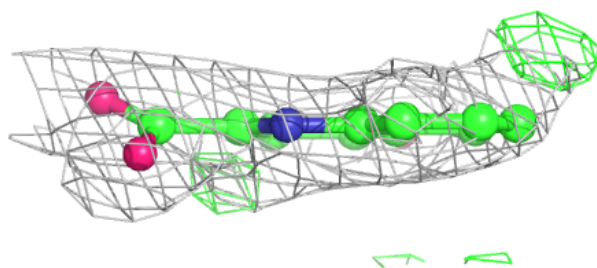
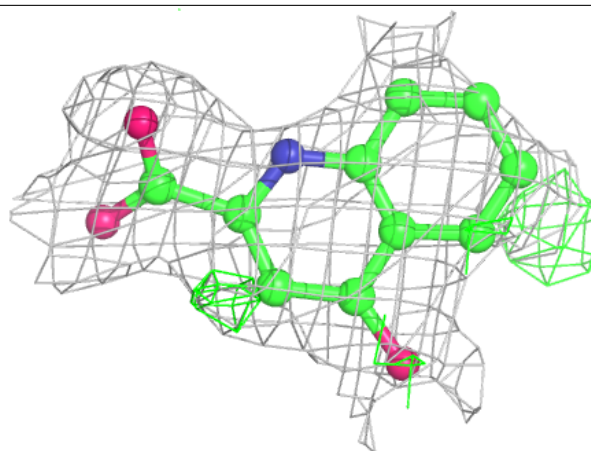
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





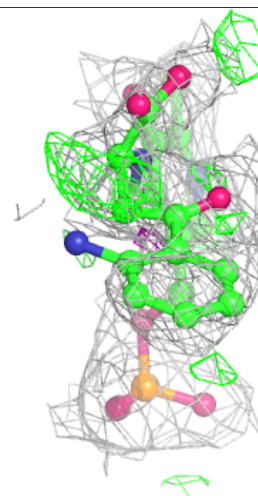
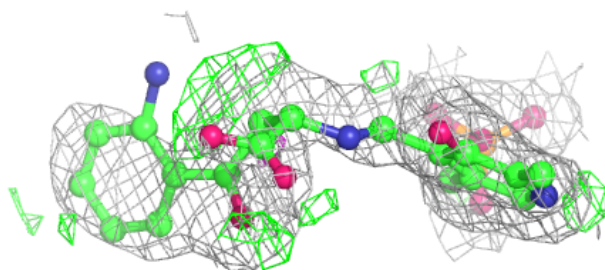
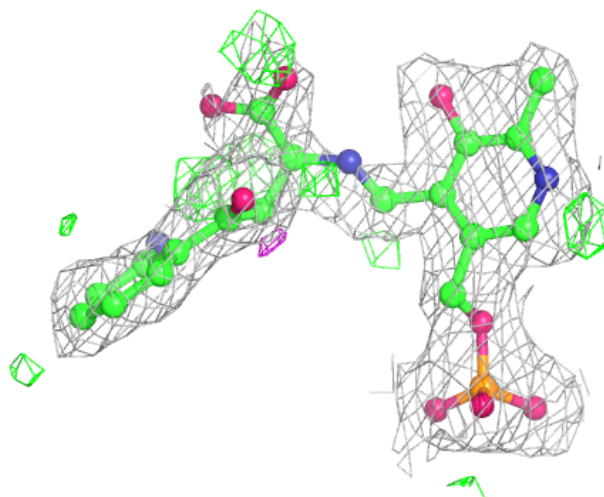
**Electron density around KYA B 501:**

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



**Electron density around Q0P A 501:**

$2mF_o-DF_c$  (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.