



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 08:32 PM UTC

PDB ID : 7LMT / pdb_00007lmt
Title : Histone-lysine N-methyltransferase NSD2-PWWP1 with compound MRT10241866a
Authors : Lei, M.; Freitas, R.F.; Dong, A.; Schapira, M.; Arrowsmith, C.H.; Edwards, A.M.; Min, J.; Structural Genomics Consortium (SGC)
Deposited on : 2021-02-05
Resolution : 2.27 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

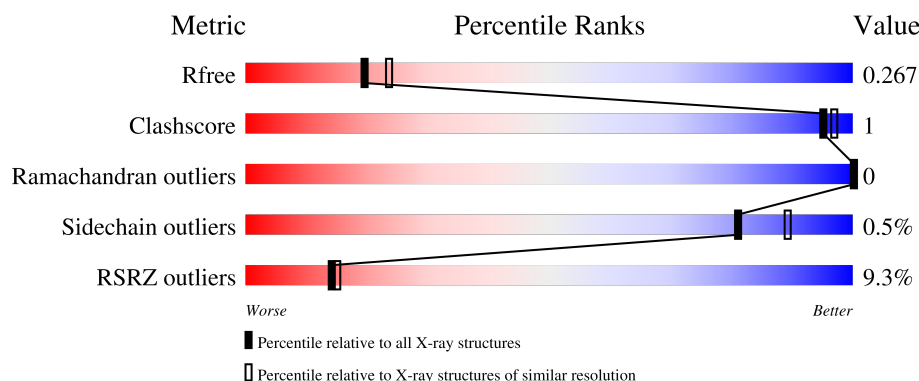
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.27 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	9078 (2.30-2.26)
Clashscore	190562	9802 (2.30-2.26)
Ramachandran outliers	187476	9690 (2.30-2.26)
Sidechain outliers	187428	9691 (2.30-2.26)
RSRZ outliers	180081	9085 (2.30-2.26)

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	140	9% 86% 6% 9%
1	B	140	7% 89% • 6%
1	C	140	11% 88% • 10%
1	D	140	9% 88% 5% 7%

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Mol	Chain	Length	Quality of chain
1	E	140	8% 84% 12%
1	F	140	10% 89% 9%
1	G	140	4% 79% 7% 14%
1	H	140	9% 86% 10%

2 Entry composition [i](#)

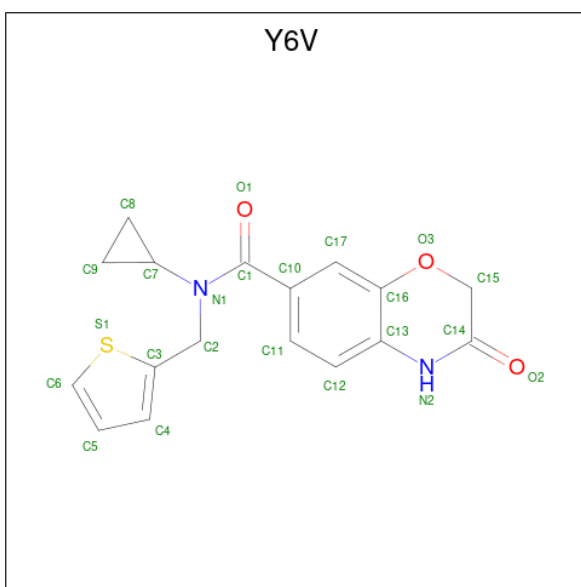
There are 4 unique types of molecules in this entry. The entry contains 7678 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Histone-lysine N-methyltransferase NSD2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	128	Total	C	N	O	S	0	0	0
			951	623	156	167	5			
1	B	131	Total	C	N	O	S	0	0	0
			963	630	158	170	5			
1	C	126	Total	C	N	O	S	0	0	0
			922	600	153	164	5			
1	D	130	Total	C	N	O	S	0	0	0
			946	618	157	166	5			
1	E	123	Total	C	N	O	S	0	0	0
			932	608	155	164	5			
1	F	127	Total	C	N	O	S	0	0	0
			931	607	155	164	5			
1	G	121	Total	C	N	O	S	0	0	0
			914	601	147	161	5			
1	H	126	Total	C	N	O	S	0	0	0
			919	599	152	163	5			

- Molecule 2 is {N}-cyclopropyl-3-oxidanylidene- {N}-(thiophen-2-ylmethyl)-4 {H}-1,4-benzoxazine-7-carboxamide (CCD ID: Y6V) (formula: C₁₇H₁₆N₂O₃S).



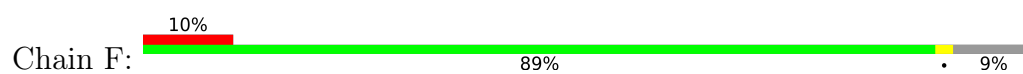
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			23	17	2	3	1		
2	B	1	Total	C	N	O	S	0	0
			23	17	2	3	1		
2	C	1	Total	C	N	O	S	0	0
			23	17	2	3	1		
2	D	1	Total	C	N	O	S	0	0
			23	17	2	3	1		
2	E	1	Total	C	N	O	S	0	0
			23	17	2	3	1		
2	F	1	Total	C	N	O	S	0	0
			23	17	2	3	1		
2	G	1	Total	C	N	O	S	0	0
			23	17	2	3	1		
2	H	1	Total	C	N	O	S	0	0
			23	17	2	3	1		

- Molecule 3 is UNKNOWN ATOM OR ION (CCD ID: UNX) (formula: X).

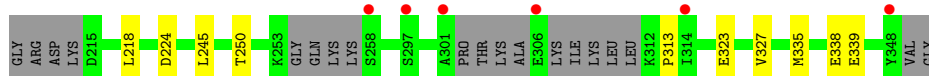
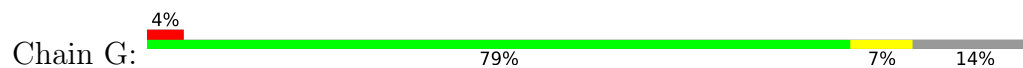
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	X	0	0
			1	1		
3	E	1	Total	X	0	0
			1	1		

- Molecule 4 is water.

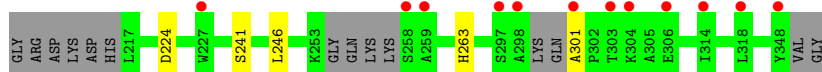
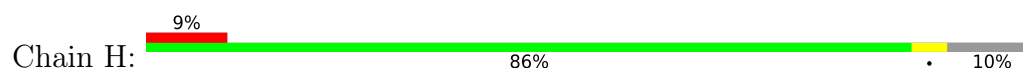
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	C	6	Total 6	O 6	0	0
4	D	3	Total 3	O 3	0	0
4	E	2	Total 2	O 2	0	0
4	F	1	Total 1	O 1	0	0
4	H	2	Total 2	O 2	0	0



- Molecule 1: Histone-lysine N-methyltransferase NSD2



- Molecule 1: Histone-lysine N-methyltransferase NSD2



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	68.30Å 69.10Å 80.19Å 75.52° 79.36° 60.47°	Depositor
Resolution (Å)	43.25 – 2.27 43.25 – 2.27	Depositor EDS
% Data completeness (in resolution range)	89.3 (43.25-2.27) 89.7 (43.25-2.27)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.47 (at 2.27Å)	Xtriage
Refinement program	BUSTER 2.10.3	Depositor
R, R_{free}	0.218 , 0.260 0.245 , 0.267	Depositor DCC
R_{free} test set	2503 reflections (4.37%)	wwPDB-VP
Wilson B-factor (Å ²)	48.8	Xtriage
Anisotropy	0.496	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 78.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.36$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	7678	wwPDB-VP
Average B, all atoms (Å ²)	70.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.02% 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: UNX, Y6V

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	0.92	0/977	1.34	6/1327 (0.5%)
1	B	0.92	0/990	1.35	3/1347 (0.2%)
1	C	0.87	0/947	1.32	0/1287
1	D	0.90	0/972	1.34	4/1323 (0.3%)
1	E	0.87	0/959	1.33	4/1301 (0.3%)
1	F	0.89	0/956	1.33	3/1300 (0.2%)
1	G	0.86	0/940	1.29	4/1276 (0.3%)
1	H	0.87	0/944	1.34	1/1283 (0.1%)
All	All	0.89	0/7685	1.33	25/10444 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1
1	H	0	1
All	All	0	2

There are no bond length outliers.

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	325	GLY	CA-C-N	6.28	128.70	120.60
1	F	325	GLY	C-N-CA	6.28	128.70	120.60
1	A	214	LYS	CA-C-N	5.71	131.70	121.66
1	A	214	LYS	C-N-CA	5.71	131.70	121.66
1	E	325	GLY	CA-C-N	5.67	127.71	120.56
1	E	325	GLY	C-N-CA	5.67	127.71	120.56

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	325	GLY	CA-C-N	5.67	127.70	120.56
1	B	325	GLY	C-N-CA	5.67	127.70	120.56
1	B	330	GLU	CB-CG-CD	5.65	122.21	112.60
1	E	259	ALA	CA-C-N	5.47	128.47	120.71
1	E	259	ALA	C-N-CA	5.47	128.47	120.71
1	H	224	ASP	CA-CB-CG	5.43	118.03	112.60
1	A	338	GLU	CB-CG-CD	5.40	121.78	112.60
1	G	338	GLU	CB-CG-CD	5.36	121.71	112.60
1	D	325	GLY	CA-C-N	5.33	127.28	120.56
1	D	325	GLY	C-N-CA	5.33	127.28	120.56
1	D	313	PRO	CA-C-N	5.26	127.66	120.35
1	D	313	PRO	C-N-CA	5.26	127.66	120.35
1	G	224	ASP	CA-CB-CG	5.24	117.84	112.60
1	F	224	ASP	CA-CB-CG	5.10	117.70	112.60
1	A	272	GLU	CB-CG-CD	5.09	121.25	112.60
1	A	325	GLY	CA-C-N	5.08	126.96	120.56
1	A	325	GLY	C-N-CA	5.08	126.96	120.56
1	G	313	PRO	CA-C-N	5.02	127.33	120.35
1	G	313	PRO	C-N-CA	5.02	127.33	120.35

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	301	ALA	Mainchain
1	H	301	ALA	Mainchain

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	951	0	840	3	0
1	B	963	0	841	3	0
1	C	922	0	793	3	0
1	D	946	0	814	4	0
1	E	932	0	829	3	0
1	F	931	0	811	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	G	914	0	796	4	0
1	H	919	0	786	2	0
2	A	23	0	0	0	0
2	B	23	0	0	0	0
2	C	23	0	0	0	0
2	D	23	0	0	1	0
2	E	23	0	0	0	0
2	F	23	0	0	0	0
2	G	23	0	0	0	0
2	H	23	0	0	0	0
3	A	1	0	0	0	0
3	E	1	0	0	0	0
4	C	6	0	0	0	0
4	D	3	0	0	0	0
4	E	2	0	0	0	0
4	F	1	0	0	0	0
4	H	2	0	0	0	0
All	All	7678	0	6510	17	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 1.

All (17) 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:246:LEU:HD11	1:D:327:VAL:HG12	1.85	0.58
1:C:335:MET:HE2	1:C:339:GLU:HB3	1.87	0.57
1:B:227:TRP:CH2	1:B:326:ILE:HD11	2.42	0.55
1:G:335:MET:HE2	1:G:339:GLU:HB3	1.89	0.54
1:A:327:VAL:HG12	1:F:246:LEU:HD11	1.90	0.53
1:B:327:VAL:HG12	1:E:246:LEU:HD11	1.93	0.49
1:G:218:LEU:HD13	1:G:250:THR:HG21	1.95	0.48
1:G:327:VAL:HG12	1:H:246:LEU:CD1	2.44	0.47
1:H:241:SER:O	1:H:263:HIS:HB3	2.15	0.47
1:D:272:GLU:HG3	2:D:401:Y6V:C17	2.47	0.44
1:A:327:VAL:HG12	1:F:246:LEU:CD1	2.47	0.43
1:B:327:VAL:HG12	1:E:246:LEU:CD1	2.49	0.43
1:E:241:SER:O	1:E:263:HIS:HB3	2.19	0.42
1:C:246:LEU:CD1	1:D:327:VAL:HG12	2.50	0.42
1:D:218:LEU:HD13	1:D:250:THR:HG21	2.02	0.41
1:A:218:LEU:HD13	1:A:250:THR:HG21	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:323:GLU:O	1:G:327:VAL:HG23	2.20	0.41

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	122/140 (87%)	121 (99%)	1 (1%)	0	100	100
1	B	127/140 (91%)	125 (98%)	2 (2%)	0	100	100
1	C	120/140 (86%)	119 (99%)	1 (1%)	0	100	100
1	D	124/140 (89%)	122 (98%)	2 (2%)	0	100	100
1	E	117/140 (84%)	117 (100%)	0	0	100	100
1	F	121/140 (86%)	121 (100%)	0	0	100	100
1	G	114/140 (81%)	112 (98%)	2 (2%)	0	100	100
1	H	120/140 (86%)	118 (98%)	2 (2%)	0	100	100
All	All	965/1120 (86%)	955 (99%)	10 (1%)	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	80/117 (68%)	79 (99%)	1 (1%)	61	76
1	B	79/117 (68%)	79 (100%)	0	100	100
1	C	75/117 (64%)	75 (100%)	0	100	100
1	D	76/117 (65%)	75 (99%)	1 (1%)	61	76
1	E	81/117 (69%)	81 (100%)	0	100	100
1	F	76/117 (65%)	76 (100%)	0	100	100
1	G	76/117 (65%)	75 (99%)	1 (1%)	61	76
1	H	73/117 (62%)	73 (100%)	0	100	100
All	All	616/936 (66%)	613 (100%)	3 (0%)	81	89

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	217	LEU
1	D	217	LEU
1	G	245	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	247	HIS
1	B	247	HIS
1	D	261	GLN
1	G	261	GLN
1	H	261	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 oligosaccharides in this entry.

5.6 Ligand geometry

Of 10 ligands modelled in this entry, 2 are unknown - leaving 8 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	Y6V	H	401	-	25,26,26	0.59	1 (4%)	34,37,37	0.95	2 (5%)
2	Y6V	E	401	-	25,26,26	0.61	1 (4%)	34,37,37	1.05	3 (8%)
2	Y6V	F	401	-	25,26,26	0.63	1 (4%)	34,37,37	0.98	3 (8%)
2	Y6V	G	401	-	25,26,26	0.56	0	34,37,37	0.93	2 (5%)
2	Y6V	D	401	-	25,26,26	0.58	1 (4%)	34,37,37	1.03	4 (11%)
2	Y6V	A	401	-	25,26,26	0.61	1 (4%)	34,37,37	0.99	3 (8%)
2	Y6V	C	401	-	25,26,26	0.67	1 (4%)	34,37,37	0.98	2 (5%)
2	Y6V	B	401	-	25,26,26	0.55	0	34,37,37	0.95	2 (5%)

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	Y6V	H	401	-	-	0/16/27/27	0/4/4/4
2	Y6V	E	401	-	-	0/16/27/27	0/4/4/4
2	Y6V	F	401	-	-	0/16/27/27	0/4/4/4
2	Y6V	G	401	-	-	0/16/27/27	0/4/4/4
2	Y6V	D	401	-	-	0/16/27/27	0/4/4/4
2	Y6V	A	401	-	-	0/16/27/27	0/4/4/4
2	Y6V	C	401	-	-	0/16/27/27	0/4/4/4
2	Y6V	B	401	-	-	0/16/27/27	0/4/4/4

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	401	Y6V	C2-N1	2.35	1.48	1.45
2	C	401	Y6V	C2-N1	2.32	1.48	1.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	401	Y6V	C2-N1	2.27	1.48	1.45
2	A	401	Y6V	C2-N1	2.20	1.48	1.45
2	H	401	Y6V	C2-N1	2.17	1.48	1.45
2	F	401	Y6V	C2-N1	2.15	1.48	1.45

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	401	Y6V	C10-C1-N1	3.07	124.80	118.72
2	E	401	Y6V	C10-C1-N1	3.00	124.66	118.72
2	E	401	Y6V	O3-C16-C13	2.94	122.75	120.74
2	A	401	Y6V	C10-C1-N1	2.88	124.42	118.72
2	C	401	Y6V	C8-C7-N1	2.84	121.99	118.46
2	G	401	Y6V	C10-C1-N1	2.74	124.15	118.72
2	F	401	Y6V	C8-C7-N1	2.74	121.86	118.46
2	B	401	Y6V	C10-C1-N1	2.60	123.86	118.72
2	H	401	Y6V	C10-C1-N1	2.54	123.75	118.72
2	C	401	Y6V	C10-C1-N1	2.50	123.67	118.72
2	F	401	Y6V	C10-C1-N1	2.48	123.63	118.72
2	F	401	Y6V	O3-C16-C13	2.44	122.41	120.74
2	H	401	Y6V	O3-C16-C13	2.30	122.31	120.74
2	A	401	Y6V	O3-C16-C13	2.27	122.29	120.74
2	D	401	Y6V	O1-C1-N1	-2.25	118.01	121.62
2	B	401	Y6V	O3-C16-C13	2.21	122.25	120.74
2	A	401	Y6V	C9-C7-N1	2.17	121.15	118.46
2	E	401	Y6V	O1-C1-N1	-2.14	118.18	121.62
2	D	401	Y6V	C9-C7-N1	2.14	121.12	118.46
2	G	401	Y6V	O3-C16-C13	2.04	122.13	120.74
2	D	401	Y6V	C13-N2-C14	-2.01	122.05	124.45

There are no chirality outliers.

There are no torsion outliers.

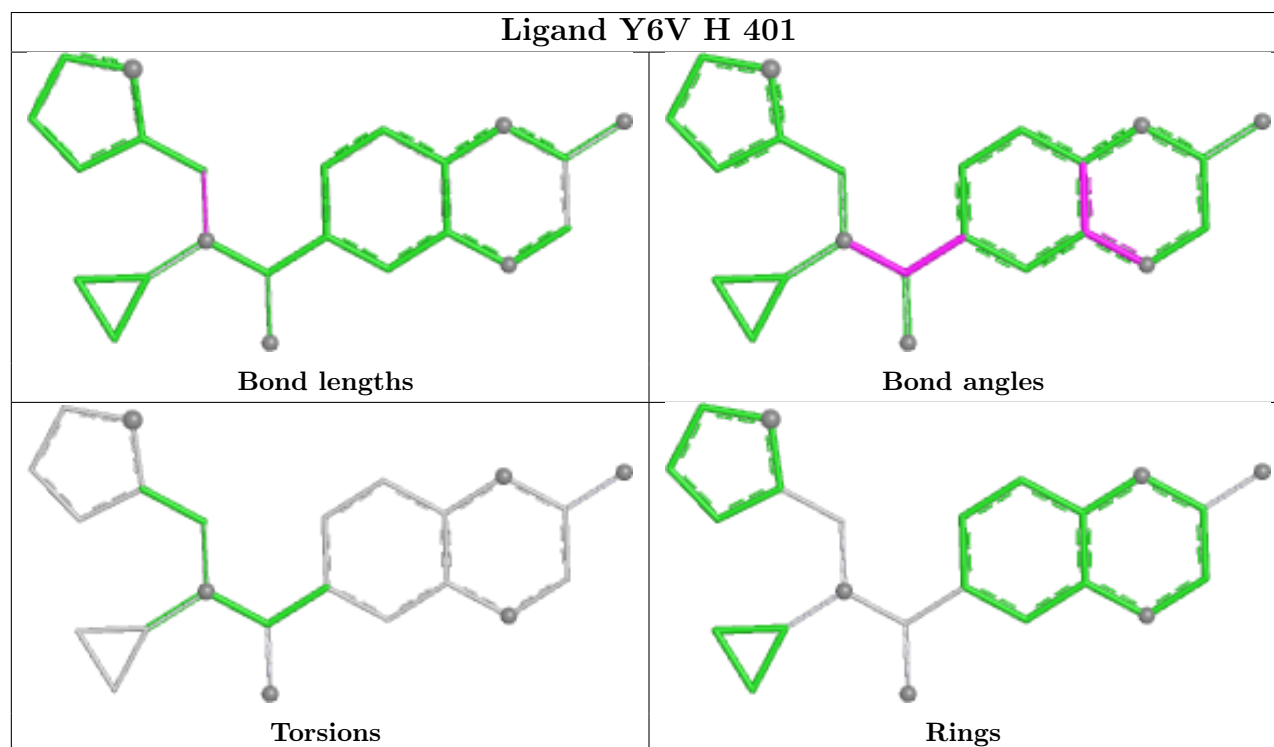
There are no ring outliers.

1 monomer is involved in 1 short contact:

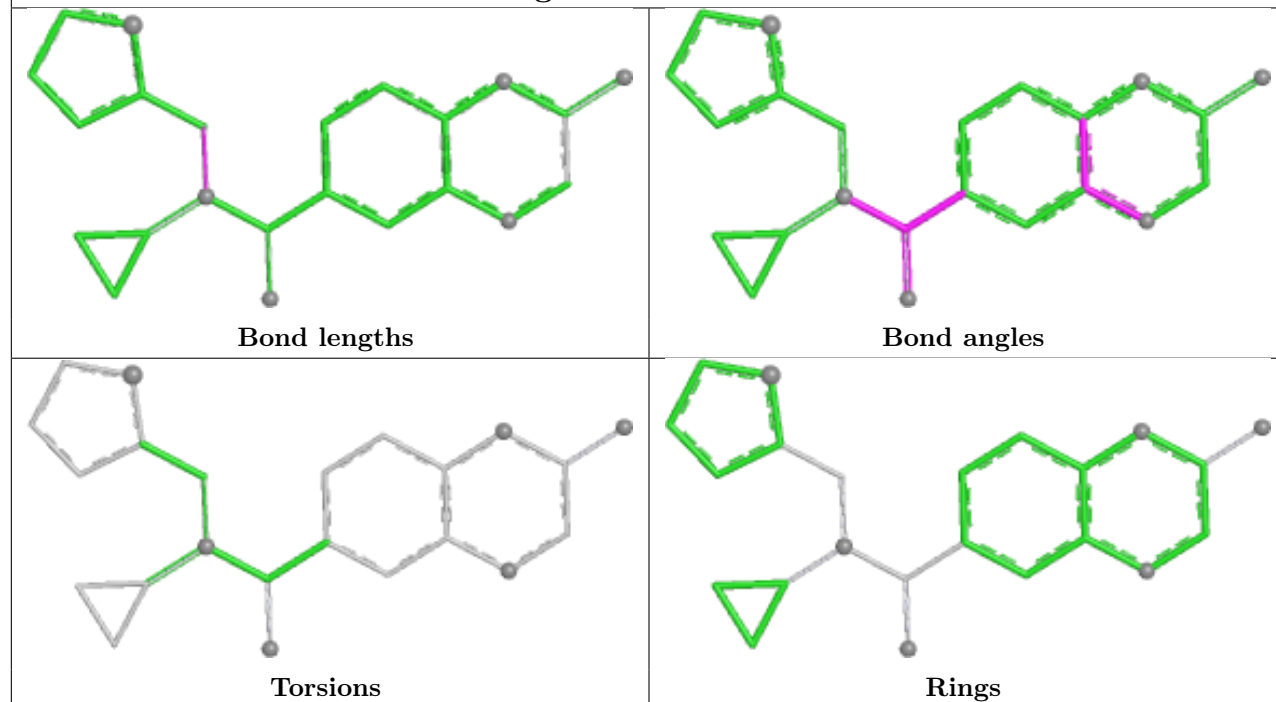
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	401	Y6V	1	0

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

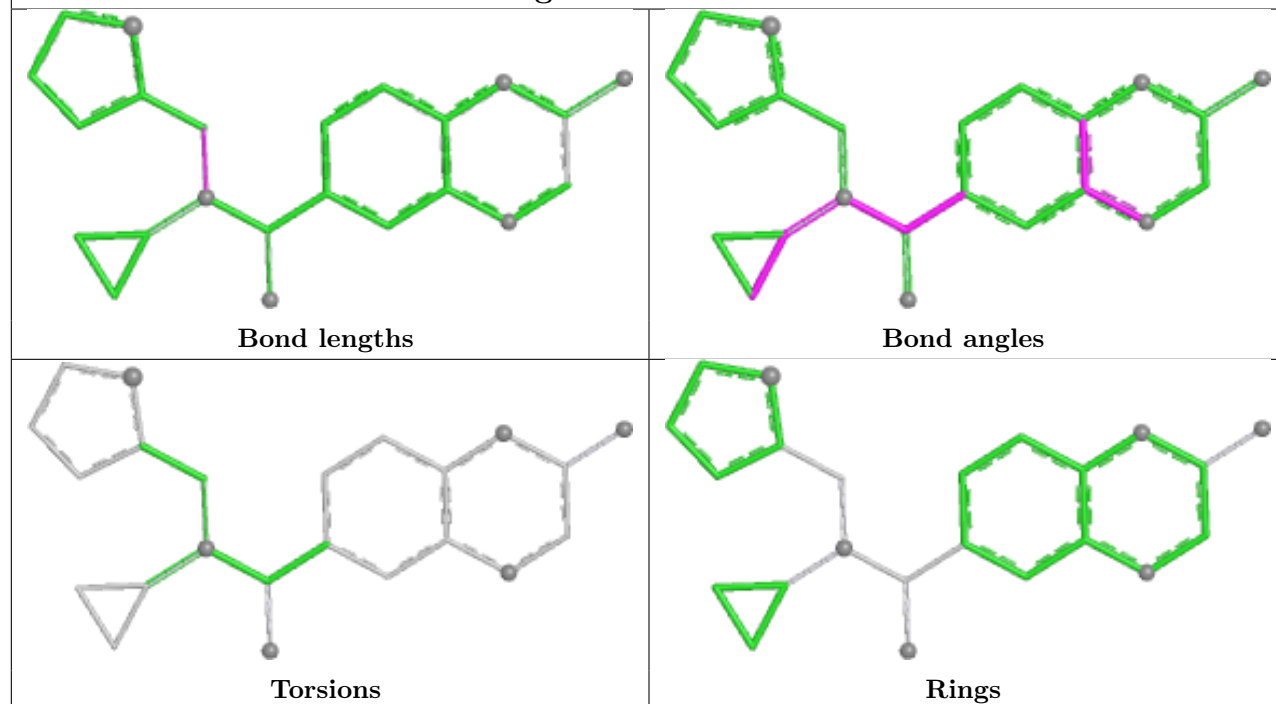
also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted 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.

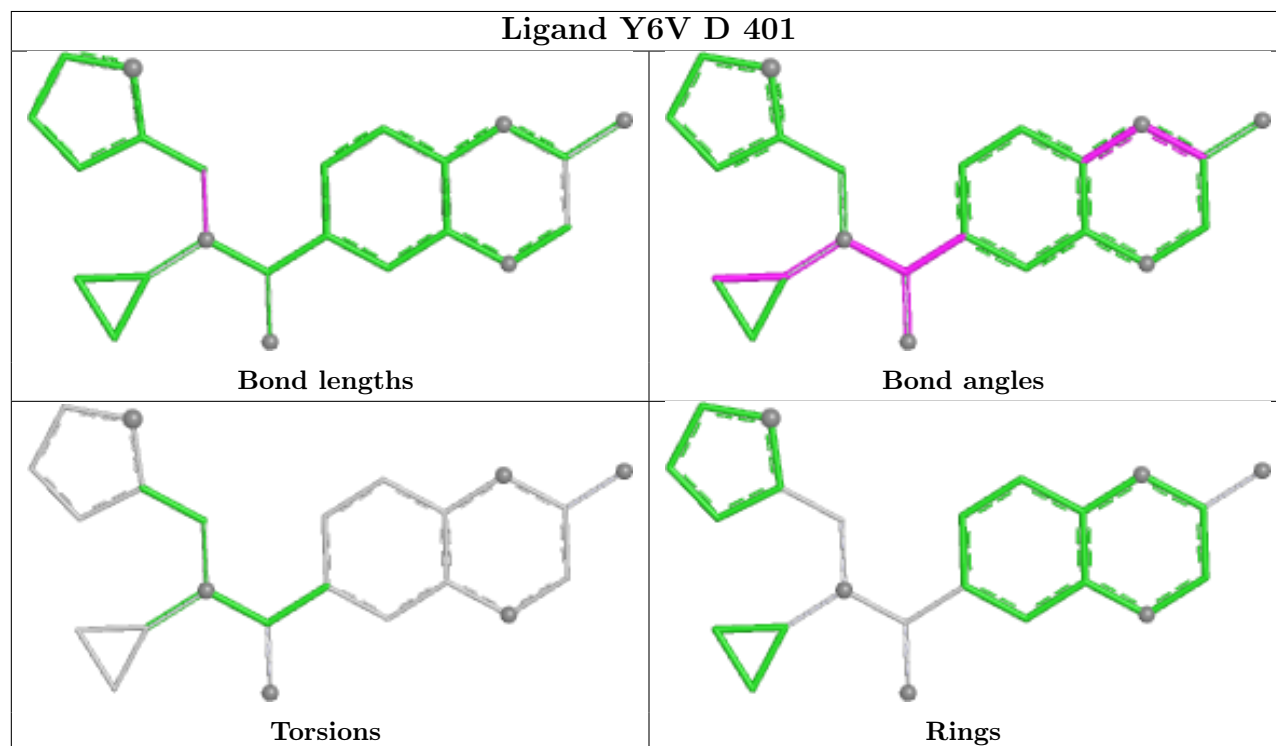
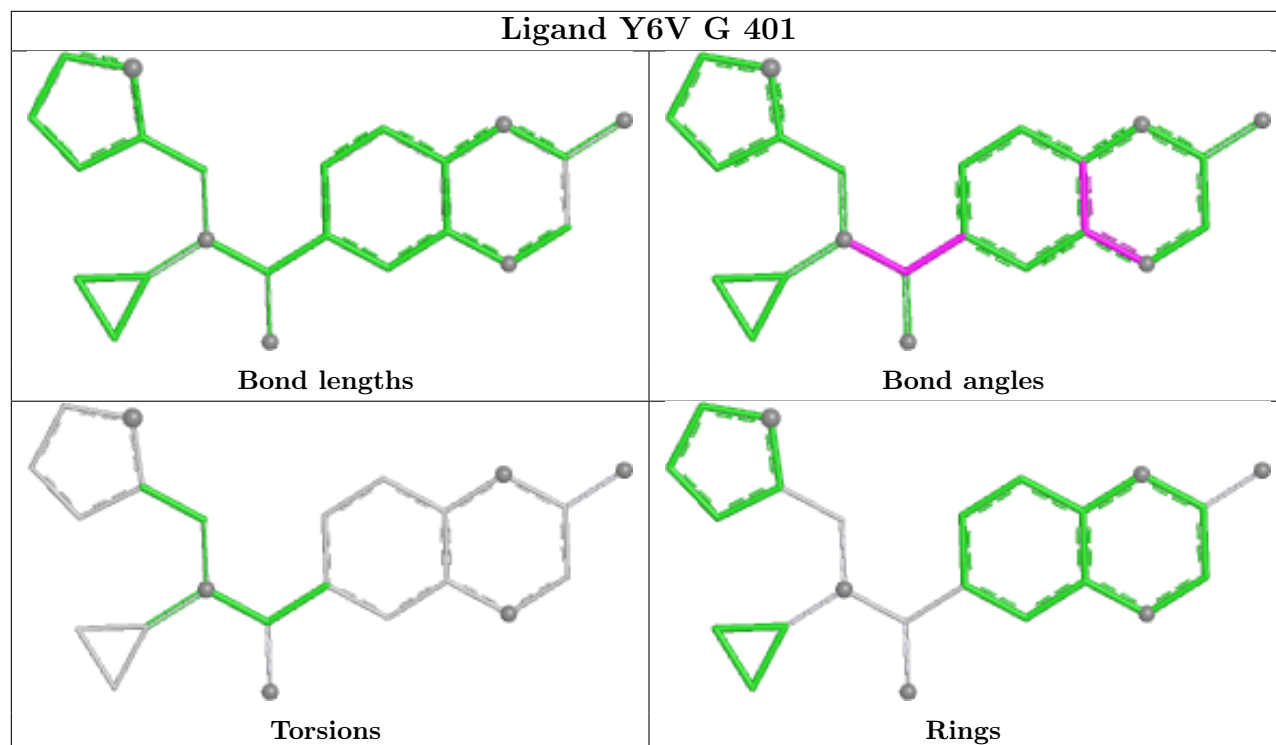


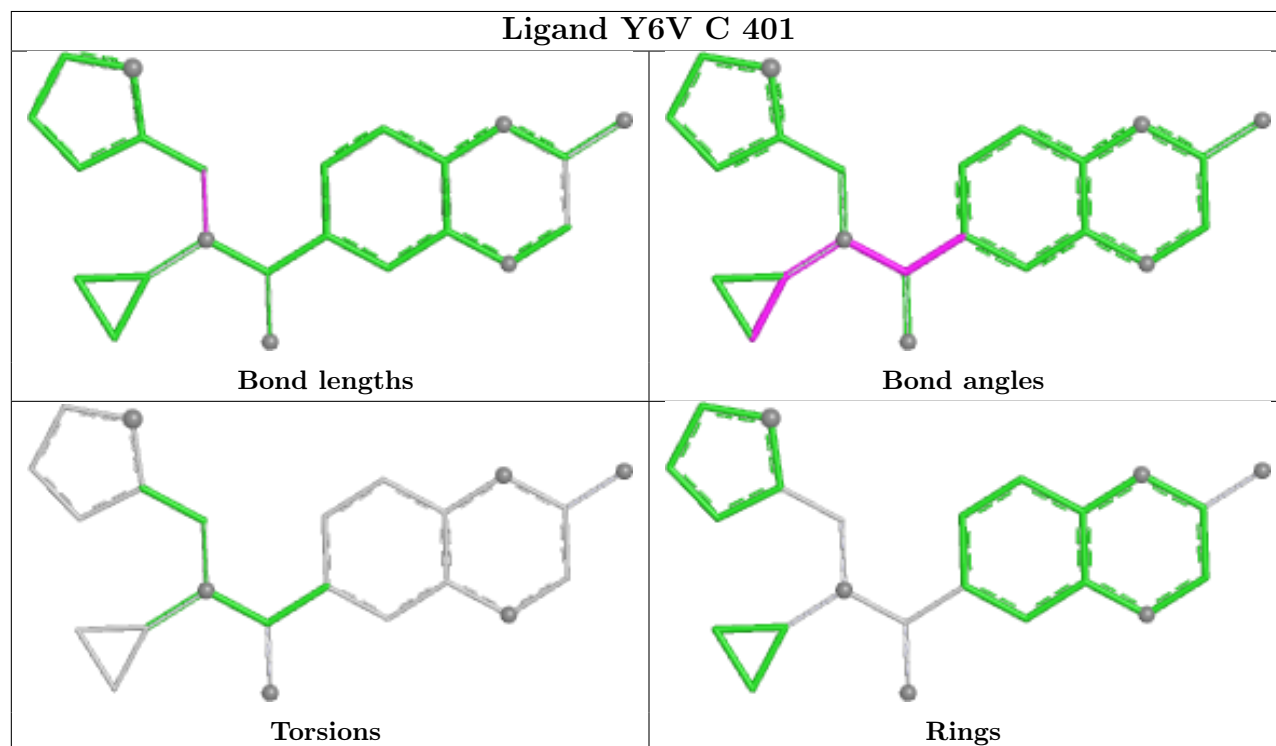
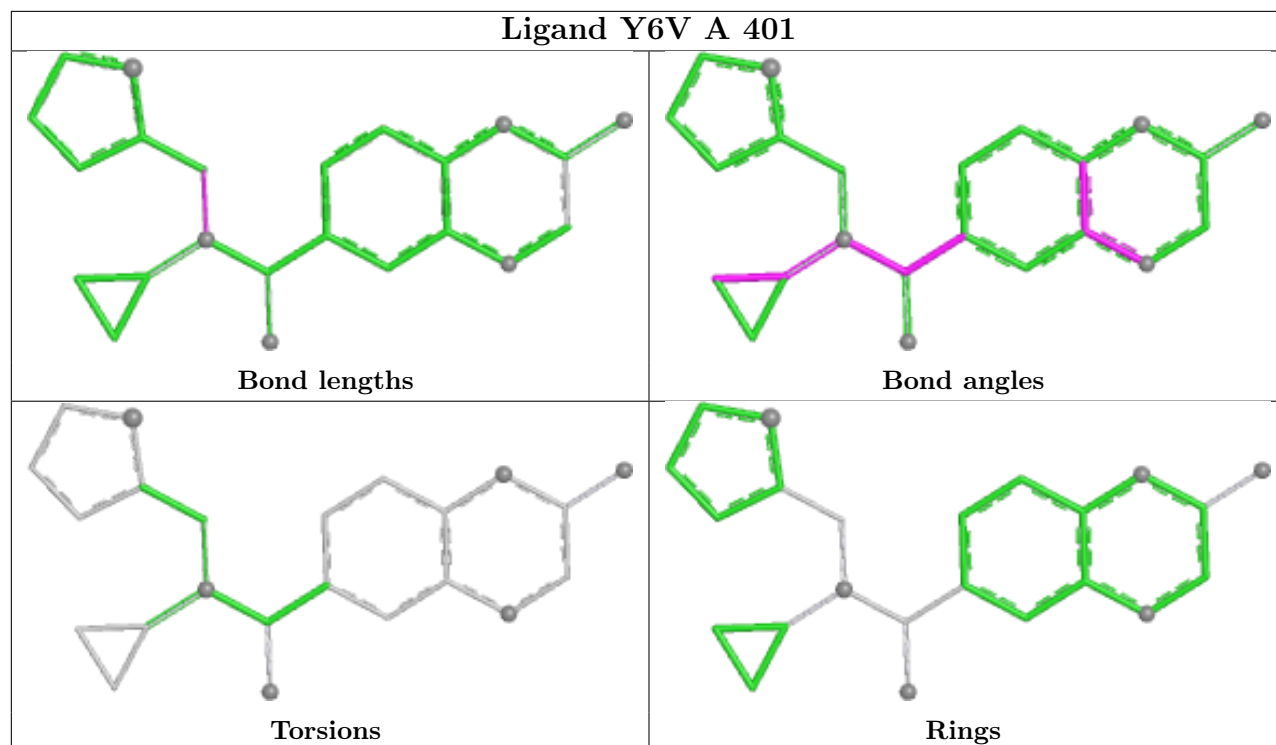
Ligand Y6V E 401

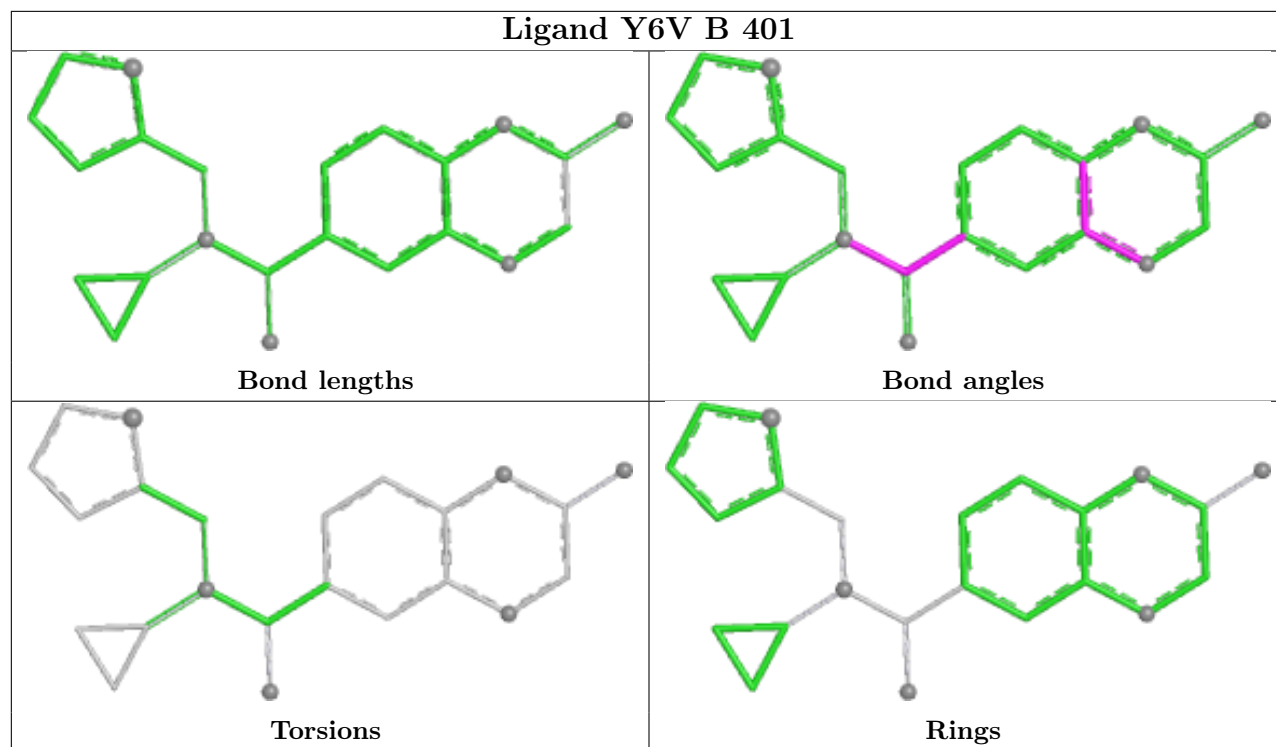


Ligand Y6V F 401









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	128/140 (91%)	0.75	13 (10%)	12 13	53, 67, 113, 137	0
1	B	131/140 (93%)	0.71	10 (7%)	20 20	52, 65, 116, 128	0
1	C	126/140 (90%)	0.76	16 (12%)	8 8	52, 67, 120, 130	0
1	D	130/140 (92%)	0.75	12 (9%)	14 15	51, 68, 114, 135	0
1	E	123/140 (87%)	0.61	11 (8%)	15 16	53, 67, 105, 117	0
1	F	127/140 (90%)	0.77	14 (11%)	10 11	49, 66, 109, 121	0
1	G	121/140 (86%)	0.64	6 (4%)	34 35	53, 67, 96, 135	0
1	H	126/140 (90%)	0.76	12 (9%)	14 15	52, 71, 130, 134	0
All	All	1012/1120 (90%)	0.72	94 (9%)	14 15	49, 67, 115, 137	0

All (94) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	348	TYR	7.2
1	F	348	TYR	5.9
1	E	314	ILE	4.5
1	A	303	THR	4.4
1	H	348	TYR	4.3
1	E	254	GLY	4.3
1	F	258	SER	4.1
1	F	303	THR	4.0
1	D	215	ASP	4.0
1	H	258	SER	3.8
1	A	214	LYS	3.7
1	C	348	TYR	3.6
1	H	306	GLU	3.6
1	E	258	SER	3.5
1	B	314	ILE	3.4
1	F	301	ALA	3.4

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Mol	Chain	Res	Type	RSRZ
1	G	314	ILE	3.3
1	E	311	LEU	3.3
1	A	348	TYR	3.2
1	A	307	LYS	3.2
1	E	310	LEU	3.2
1	E	348	TYR	3.2
1	B	214	LYS	3.1
1	H	297	SER	3.1
1	C	303	THR	3.1
1	A	258	SER	3.0
1	H	227	TRP	3.0
1	C	305	ALA	2.9
1	G	348	TYR	2.9
1	D	258	SER	2.9
1	F	314	ILE	2.9
1	B	215	ASP	2.9
1	C	258	SER	2.9
1	C	306	GLU	2.9
1	B	310	LEU	2.9
1	H	259	ALA	2.9
1	D	314	ILE	2.9
1	A	215	ASP	2.9
1	D	288	GLY	2.8
1	B	258	SER	2.8
1	G	297	SER	2.8
1	F	346	PHE	2.8
1	H	298	ALA	2.8
1	D	295	GLN	2.8
1	C	314	ILE	2.8
1	A	306	GLU	2.8
1	B	348	TYR	2.8
1	A	314	ILE	2.7
1	G	258	SER	2.7
1	F	259	ALA	2.7
1	E	346	PHE	2.7
1	G	306	GLU	2.7
1	B	306	GLU	2.6
1	G	301	ALA	2.6
1	C	307	LYS	2.6
1	F	305	ALA	2.6
1	H	301	ALA	2.6
1	D	306	GLU	2.5

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Mol	Chain	Res	Type	RSRZ
1	H	314	ILE	2.5
1	B	288	GLY	2.5
1	A	295	GLN	2.5
1	H	303	THR	2.5
1	C	288	GLY	2.5
1	A	310	LEU	2.4
1	D	303	THR	2.4
1	E	347	LEU	2.4
1	F	308	ILE	2.4
1	C	253	LYS	2.4
1	F	297	SER	2.4
1	F	306	GLU	2.4
1	C	308	ILE	2.3
1	D	299	LYS	2.3
1	F	315	SER	2.3
1	C	290	PHE	2.3
1	H	304	LYS	2.3
1	D	304	LYS	2.3
1	C	294	CYS	2.2
1	C	298	ALA	2.2
1	E	227	TRP	2.2
1	B	308	ILE	2.2
1	D	300	GLN	2.2
1	C	232	GLY	2.1
1	D	297	SER	2.1
1	C	304	LYS	2.1
1	F	307	LYS	2.1
1	A	313	PRO	2.1
1	A	342	ALA	2.1
1	E	309	LYS	2.1
1	H	318	LEU	2.1
1	C	230	VAL	2.1
1	A	299	LYS	2.0
1	B	347	LEU	2.0
1	F	277	PHE	2.0
1	E	293	LEU	2.0

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

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](#)

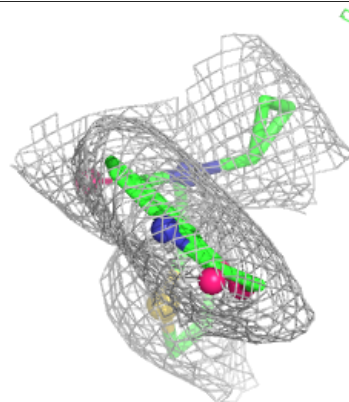
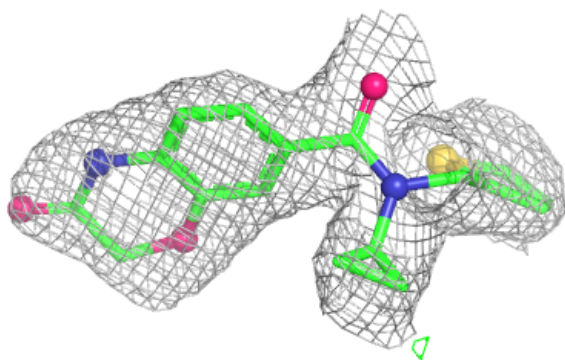
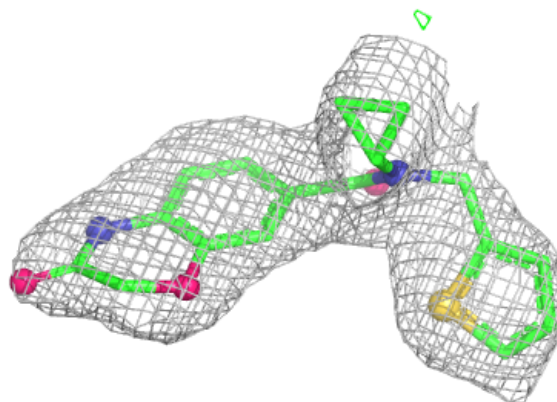
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	UNX	A	402	1/1	0.81	0.13	44,44,44,44	0
3	UNX	E	402	1/1	0.85	0.19	41,41,41,41	0
2	Y6V	F	401	23/23	0.94	0.09	48,59,68,69	0
2	Y6V	A	401	23/23	0.94	0.09	55,62,66,66	0
2	Y6V	E	401	23/23	0.94	0.08	52,57,64,68	0
2	Y6V	G	401	23/23	0.95	0.08	56,62,69,72	0
2	Y6V	B	401	23/23	0.95	0.08	52,59,62,66	0
2	Y6V	D	401	23/23	0.95	0.09	57,61,62,62	0
2	Y6V	C	401	23/23	0.96	0.07	51,56,64,66	0
2	Y6V	H	401	23/23	0.96	0.08	60,65,67,71	0

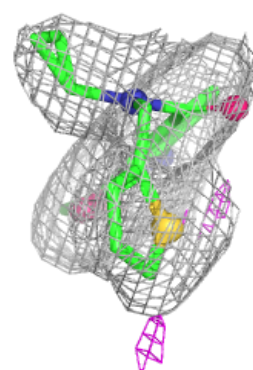
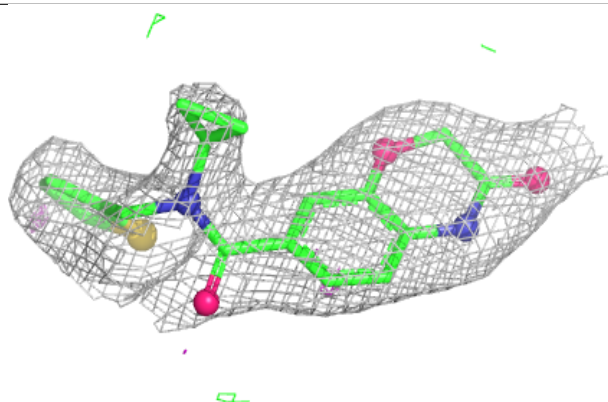
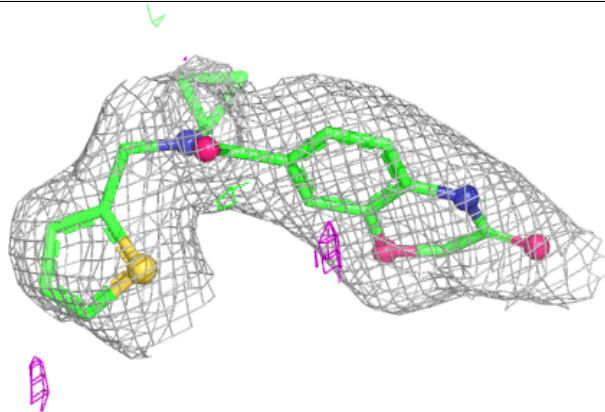
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around Y6V F 401:

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

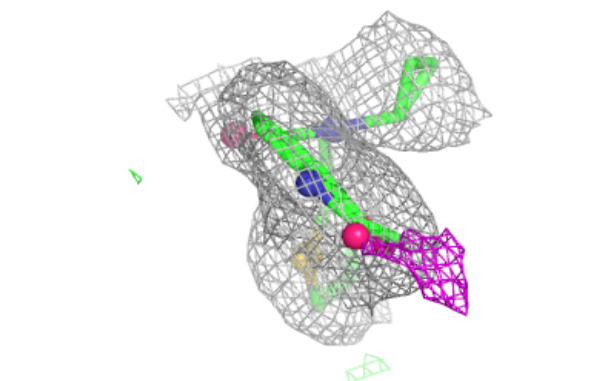
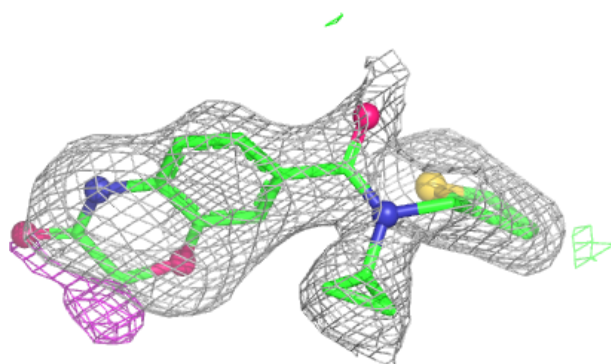
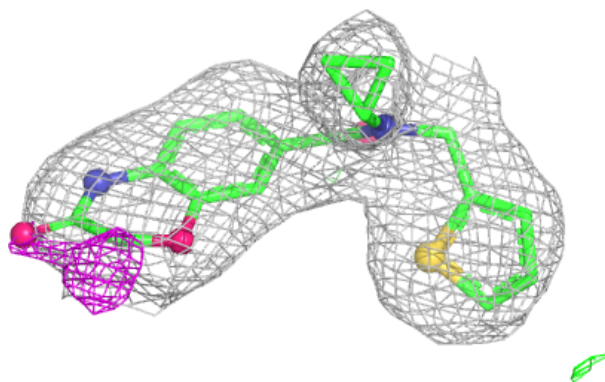
**Electron density around Y6V A 401:**

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

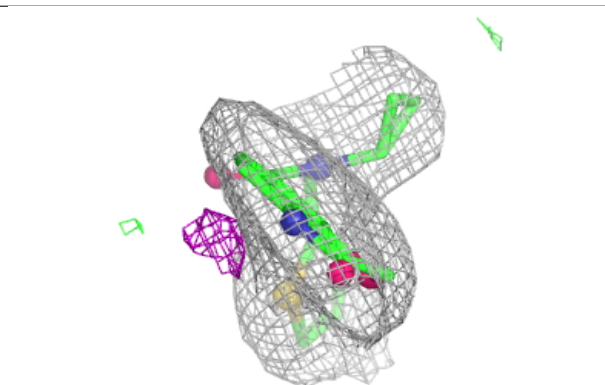
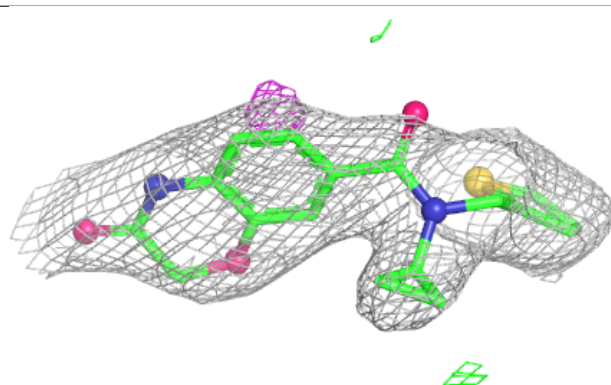
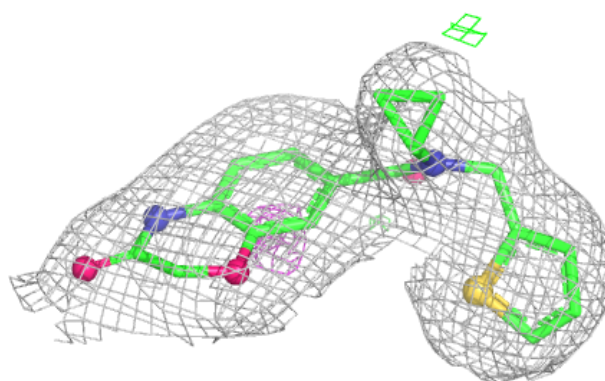


Electron density around Y6V E 401:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

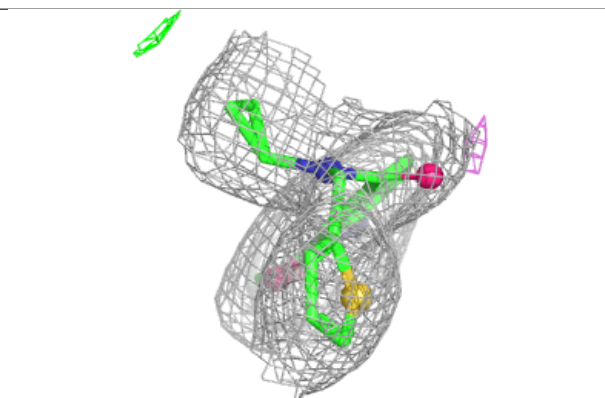
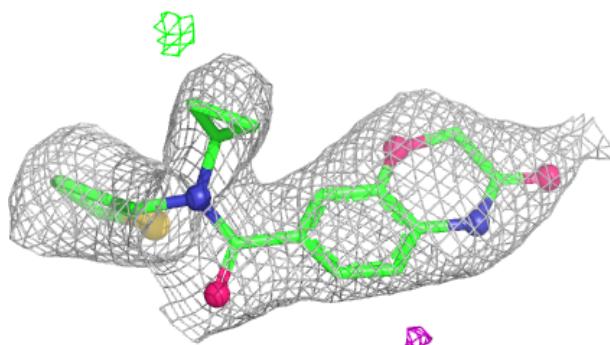
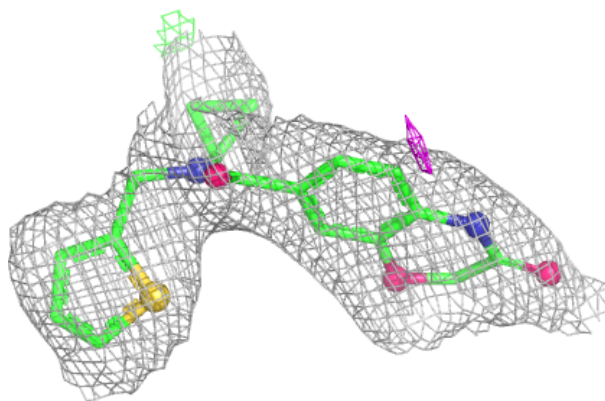
**Electron density around Y6V G 401:**

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and green (positive)

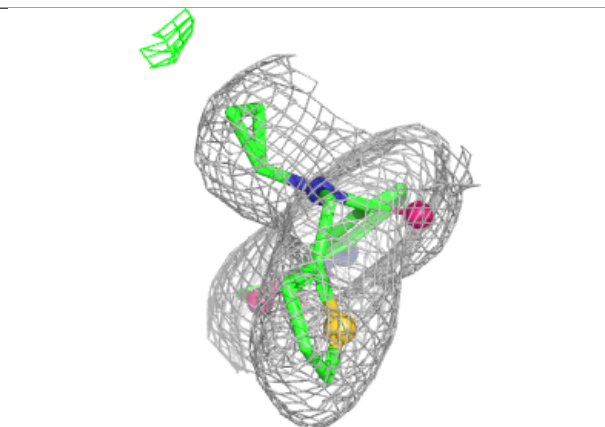
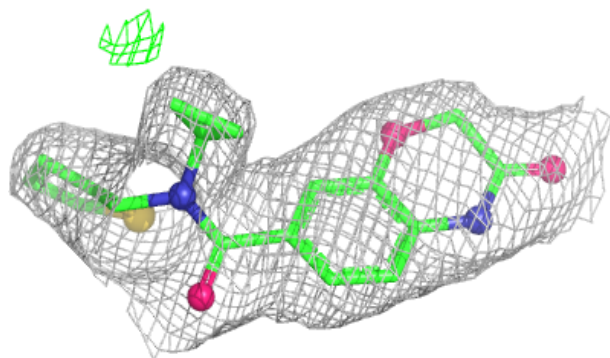
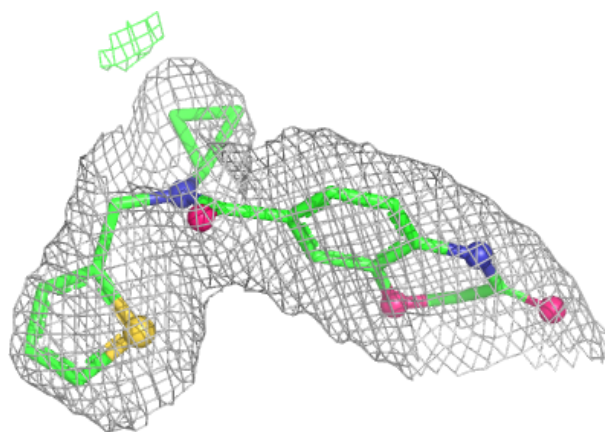


Electron density around Y6V B 401:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

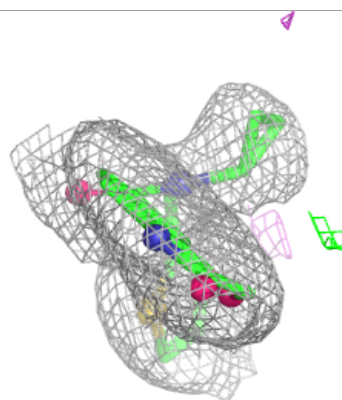
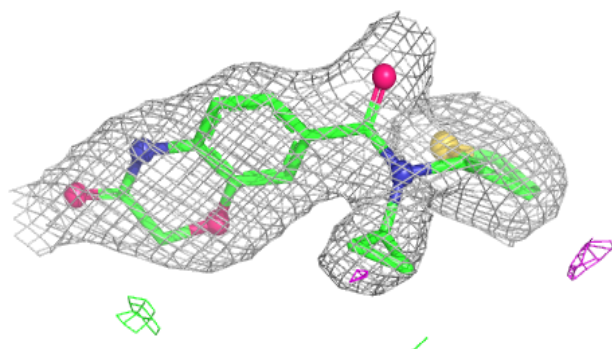
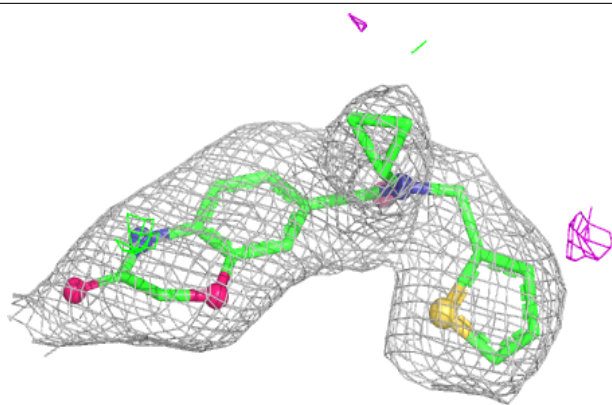
**Electron density around Y6V D 401:**

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

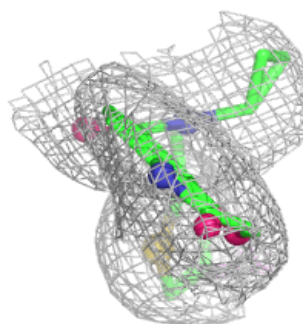
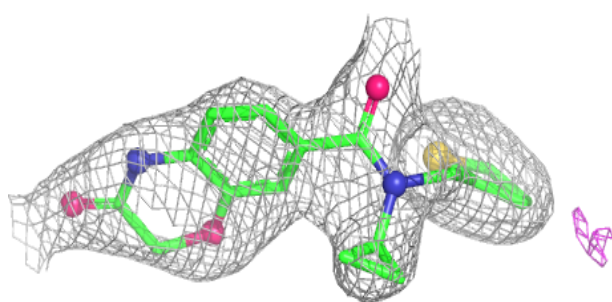
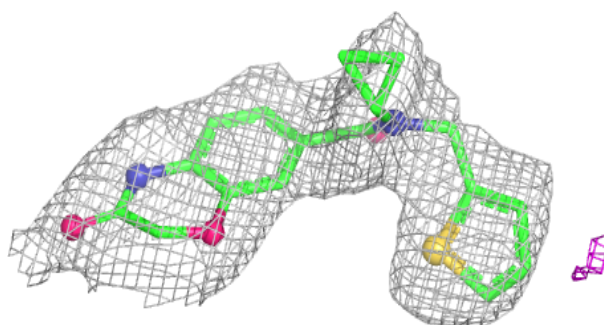


Electron density around Y6V C 401:

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

**Electron density around Y6V H 401:**

$2mF_o-DF_c$ (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.