



Full wwPDB X-ray Structure Validation Report ⓘ

May 13, 2020 – 03:21 am BST

PDB ID : 5IXS
Title : Lactate Dehydrogenase in complex with hydroxylactam inhibitor compound
9: (6R)-3-[(2-chlorophenyl)sulfanyl]-4-hydroxy-6-(3-hydroxyphenyl)-6-(thiophen-3-yl)-5,6-dihydropyridin-2(1H)-one
Authors : Ultsch, M.; Eigenbrot, C.
Deposited on : 2016-03-23
Resolution : 2.05 Å(reported)

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

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

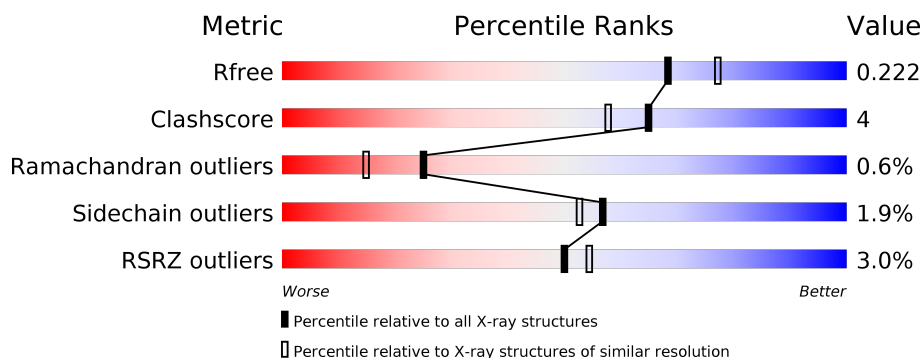
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.05 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	1692 (2.04-2.04)
Clashscore	141614	1773 (2.04-2.04)
Ramachandran outliers	138981	1752 (2.04-2.04)
Sidechain outliers	138945	1752 (2.04-2.04)
RSRZ outliers	127900	1672 (2.04-2.04)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	331	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 1%, green 98%);"></div> <div style="display: flex; justify-content: space-between; margin-top: 2px;"> % 89% 9% .. </div> </div>
1	B	331	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 2%, orange 2%, yellow 1%, green 95%);"></div> <div style="display: flex; justify-content: space-between; margin-top: 2px;"> 2% 86% 11% • </div> </div>
1	C	331	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 6%, orange 6%, yellow 1%, green 87%);"></div> <div style="display: flex; justify-content: space-between; margin-top: 2px;"> 6% 90% 7% .. </div> </div>
1	D	331	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 3%, orange 3%, yellow 1%, green 93%);"></div> <div style="display: flex; justify-content: space-between; margin-top: 2px;"> 3% 93% 5% • </div> </div>

2 Entry composition [i](#)

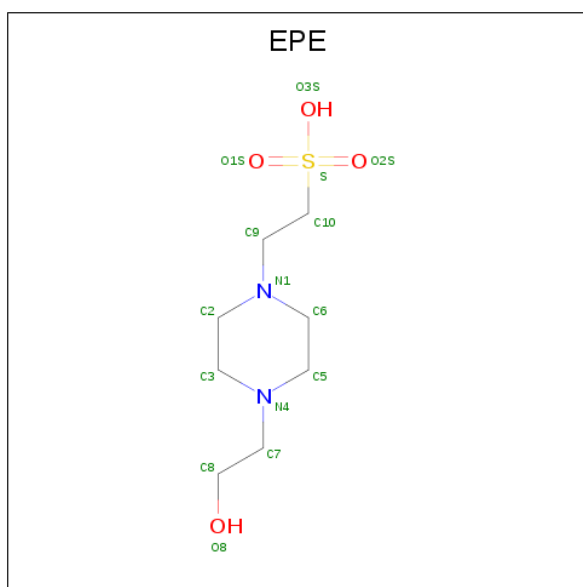
There are 7 unique types of molecules in this entry. The entry contains 11112 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 L-lactate dehydrogenase A chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	327	Total	C	N	O	S	0	4	0
			2556	1635	434	473	14			
1	B	324	Total	C	N	O	S	0	3	0
			2532	1626	430	463	13			
1	C	324	Total	C	N	O	S	0	2	0
			2522	1616	428	465	13			
1	D	330	Total	C	N	O	S	0	4	0
			2580	1650	441	475	14			

- Molecule 2 is 4-(2-HYDROXYETHYL)-1-PIPERAZINE ETHANESULFONIC ACID (three-letter code: EPE) (formula: C₈H₁₈N₂O₄S).



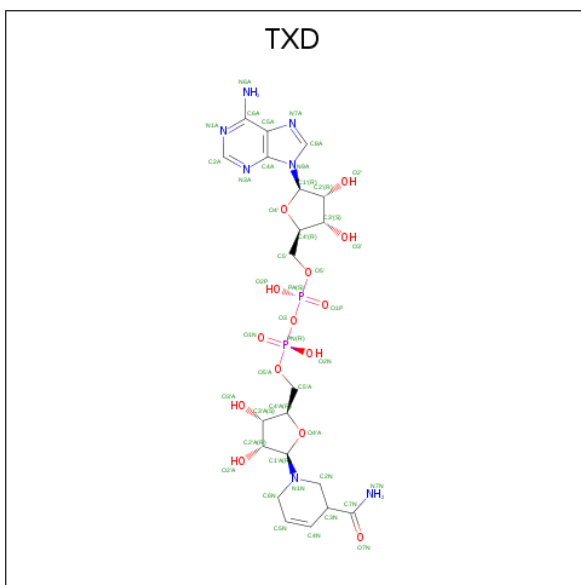
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			15	8	2	4	1		
2	B	1	Total	C	N	O	S	0	0
			15	8	2	4	1		

Continued on next page...

Continued from previous page...

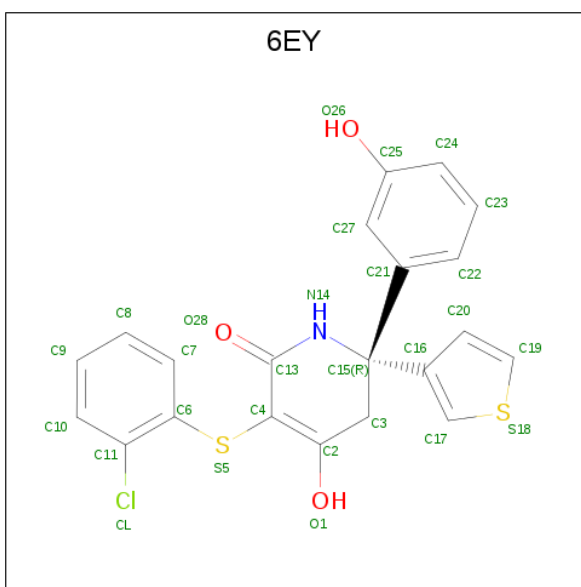
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	C	1	Total	C	N	O	S	0	0
			15	8	2	4	1		

- Molecule 3 is 1,4,5,6-TETRAHYDRONICOTINAMIDE ADENINE DINUCLEOTIDE (three-letter code: TXD) (formula: $C_{21}H_{31}N_7O_{14}P_2$).



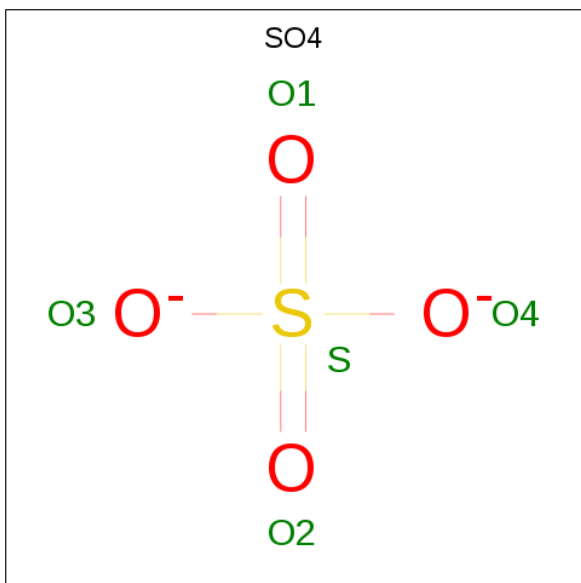
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total 44	C 21	N 7	O 14	P 2	0	0
3	B	1	Total 44	C 21	N 7	O 14	P 2	0	0
3	D	1	Total 44	C 21	N 7	O 14	P 2	0	0

- Molecule 4 is (6R)-3-[(2-chlorophenyl)sulfanyl]-4-hydroxy-6-(3-hydroxyphenyl)-6-(thiophen-3-yl)-5,6-dihydropyridin-2(1H)-one (three-letter code: 6EY) (formula: $C_{21}H_{16}ClNO_3S_2$).



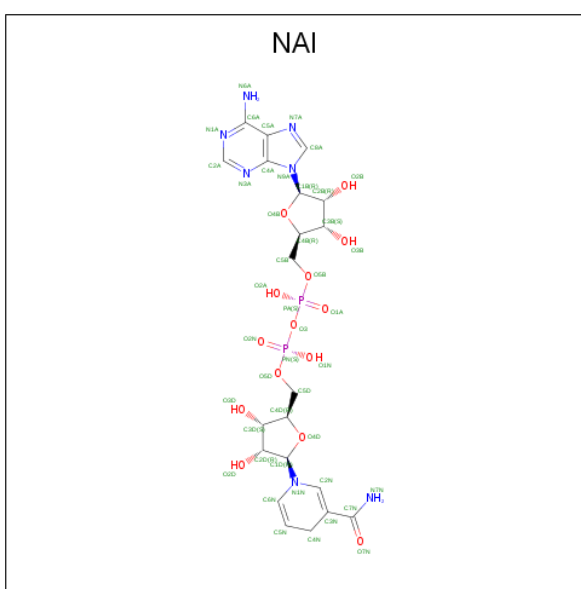
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
4	A	1	Total 28	C 21	Cl 1	N 1	O 3	S 2	0	0
4	B	1	Total 28	C 21	Cl 1	N 1	O 3	S 2	0	0
4	C	1	Total 28	C 21	Cl 1	N 1	O 3	S 2	0	0
4	D	1	Total 28	C 21	Cl 1	N 1	O 3	S 2	0	0

- Molecule 5 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total O S 5 4 1	0	0
5	B	1	Total O S 5 4 1	0	0
5	C	1	Total O S 5 4 1	0	0
5	D	1	Total O S 5 4 1	0	0

- Molecule 6 is 1,4-DIHYDRONICOTINAMIDE ADENINE DINUCLEOTIDE (three-letter code: NAI) (formula: $C_{21}H_{29}N_7O_{14}P_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	C	1	Total	C	N	O	P	0	0
			44	21	7	14	2		

- Molecule 7 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	158	Total O 158 158	0	0
7	B	159	Total O 159 159	0	0
7	C	98	Total O 98 98	0	0
7	D	154	Total O 154 154	0	0

- Molecule 1: L-lactate dehydrogenase A chain





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	74.62Å 81.23Å 104.80Å 90.00° 97.56° 90.00°	Depositor
Resolution (Å)	37.80 – 2.05 37.83 – 2.05	Depositor EDS
% Data completeness (in resolution range)	98.0 (37.80-2.05) 98.1 (37.83-2.05)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.71 (at 2.05Å)	Xtriage
Refinement program	REFMAC 5.8.0049	Depositor
R, R_{free}	0.174 , 0.221 0.178 , 0.222	Depositor DCC
R_{free} test set	1571 reflections (2.05%)	wwPDB-VP
Wilson B-factor (Å ²)	32.7	Xtriage
Anisotropy	0.020	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 36.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	11112	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.97% 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: NAI, TXD, EPE, SO4, 6EY

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	0.40	0/2611	0.62	1/3529 (0.0%)
1	B	0.39	0/2585	0.58	1/3495 (0.0%)
1	C	0.37	0/2571	0.56	0/3476
1	D	0.38	0/2635	0.56	0/3560
All	All	0.39	0/10402	0.58	2/14060 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	105	ARG	NE-CZ-NH1	6.40	123.50	120.30
1	B	302	LEU	CA-CB-CG	5.75	128.53	115.30

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	2556	0	2656	32	0
1	B	2532	0	2639	29	0
1	C	2522	0	2623	18	0
1	D	2580	0	2687	16	0
2	A	15	0	17	0	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	15	0	17	1	0
2	C	15	0	18	0	0
3	A	44	0	27	6	0
3	B	44	0	27	6	0
3	D	44	0	27	2	0
4	A	28	0	0	2	0
4	B	28	0	0	2	0
4	C	28	0	0	0	0
4	D	28	0	0	0	0
5	A	5	0	0	0	0
5	B	5	0	0	0	0
5	C	5	0	0	0	0
5	D	5	0	0	0	0
6	C	44	0	26	1	0
7	A	158	0	0	2	0
7	B	159	0	0	5	0
7	C	98	0	0	1	0
7	D	154	0	0	0	0
All	All	11112	0	10764	87	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 4.

All (87) 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:105:ARG:HG2	1:A:105:ARG:HH11	1.23	0.97
1:A:20:ASN:OD1	7:A:501:HOH:O	2.01	0.78
1:B:98:ARG:HG3	4:B:403:6EY:C24	2.16	0.76
1:B:276:ILE:HD11	1:B:286:VAL:HG23	1.66	0.76
1:A:14:GLU:HB3	1:A:15:GLU:HB2	1.69	0.74
1:B:203:MET:HB2	7:B:654:HOH:O	1.91	0.70
1:B:30:VAL:HG21	3:B:402:TXD:H6N1	1.71	0.70
1:A:105:ARG:HG2	1:A:105:ARG:NH1	1.93	0.69
1:B:203:MET:SD	7:B:654:HOH:O	2.52	0.67
1:A:247:THR:HG23	3:A:402:TXD:H5N	1.78	0.66
1:A:293:ILE:CD1	1:C:179:VAL:HG22	2.25	0.65
1:B:10:ASN:OD1	1:B:13:LYS:HG2	1.97	0.64
1:A:99:GLN:HG2	1:A:108:LEU:HD22	1.82	0.62
1:B:72[A]:ARG:NH1	7:B:502:HOH:O	2.32	0.61
1:A:247:THR:CG2	3:A:402:TXD:H5N	2.30	0.61

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:96:GLY:O	1:D:111:ARG:NH1	2.36	0.59
1:A:3:LEU:HD13	1:B:214:HIS:HB2	1.84	0.59
1:D:203[A]:MET:HE2	1:D:217:LEU:HD21	1.86	0.58
1:D:203[A]:MET:HG2	1:D:210:LEU:HD22	1.85	0.58
1:B:137:ASN:ND2	3:B:402:TXD:H2N1	2.18	0.58
1:A:105:ARG:HG3	1:A:106:LEU:N	2.19	0.57
1:C:13:LYS:O	1:C:15:GLU:N	2.37	0.57
4:A:403:6EY:O26	7:A:502:HOH:O	2.17	0.56
1:B:98:ARG:CG	4:B:403:6EY:C24	2.84	0.56
1:B:276:ILE:HD13	1:B:282:ILE:HD13	1.87	0.56
1:B:225:GLN:NE2	7:B:503:HOH:O	2.38	0.56
1:C:276:ILE:HD11	1:C:286:VAL:HG23	1.88	0.56
1:A:293:ILE:HD12	1:C:179:VAL:HG22	1.87	0.55
1:A:105:ARG:CG	1:A:105:ARG:HH11	2.08	0.55
1:A:14:GLU:HB3	1:A:15:GLU:CB	2.35	0.55
1:A:204:ASN:HA	1:A:210:LEU:HD13	1.90	0.54
1:A:30:VAL:HG21	3:A:402:TXD:H6N1	1.89	0.54
1:D:30:VAL:HG21	3:D:801:TXD:H6N1	1.89	0.53
1:C:154[B]:LYS:HE2	1:C:275:MET:SD	2.49	0.52
1:C:182:LEU:HD11	1:D:72[B]:ARG:NH2	2.24	0.52
1:D:203[A]:MET:CE	1:D:217:LEU:HD21	2.40	0.52
1:D:115:ILE:HG22	1:D:119:ILE:HD12	1.93	0.51
1:D:200:TRP:CE3	1:D:203[B]:MET:SD	3.04	0.51
1:A:104:SER:O	1:A:105:ARG:HB3	2.11	0.50
1:A:293:ILE:HG13	1:A:301:ASP:HB2	1.93	0.50
1:A:104:SER:HB2	1:A:107:ASN:HB2	1.92	0.50
1:A:97:ALA:HA	3:A:402:TXD:O2P	2.12	0.50
1:C:115:ILE:HG23	6:C:402:NAI:N6A	2.27	0.49
1:A:96:GLY:O	1:A:97:ALA:CB	2.61	0.49
1:A:96:GLY:O	1:A:97:ALA:HB3	2.13	0.49
1:B:153:PRO:HA	2:B:401:EPE:H92	1.96	0.48
1:A:115:ILE:HG23	3:A:402:TXD:N6A	2.29	0.48
1:D:97:ALA:HB3	1:D:111:ARG:HH12	1.80	0.47
1:D:293[A]:ILE:HD12	1:D:301:ASP:HB2	1.96	0.47
1:A:82:TYR:CG	1:A:122:ASN:HB3	2.49	0.47
1:A:203:MET:HG2	1:A:210:LEU:HD22	1.96	0.47
1:D:137:ASN:ND2	3:D:801:TXD:H2N1	2.29	0.47
1:C:182:LEU:HD11	1:D:72[B]:ARG:CZ	2.46	0.46
1:C:197:VAL:HG21	1:C:315:LEU:HD23	1.96	0.46
1:B:109:VAL:O	1:B:113:VAL:HG23	2.15	0.46
1:D:12:LEU:HD22	1:D:14:GLU:HG2	1.96	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:154[B]:LYS:HE3	7:C:515:HOH:O	2.15	0.46
1:B:293:ILE:HD12	1:B:301:ASP:HB2	1.98	0.46
1:B:270:HIS:CD2	1:B:294:LEU:HD12	2.52	0.45
1:B:98:ARG:NH2	3:B:402:TXD:O1P	2.49	0.45
1:D:108:LEU:HD12	1:D:111:ARG:HH21	1.82	0.45
1:D:241:ILE:CD1	1:D:247:THR:HG23	2.47	0.45
1:A:98:ARG:O	1:A:99:GLN:C	2.54	0.45
1:B:16:GLN:HE21	1:C:297:ASN:HD22	1.65	0.44
1:B:115:ILE:HG23	3:B:402:TXD:N6A	2.33	0.44
1:A:98:ARG:HG2	4:A:403:6EY:O26	2.17	0.44
1:A:137:ASN:ND2	3:A:402:TXD:H2N1	2.33	0.43
1:B:82:TYR:CG	1:B:122:ASN:HB3	2.53	0.43
1:A:14:GLU:HB3	1:A:15:GLU:CA	2.48	0.43
1:B:231:LYS:HA	1:B:231:LYS:HD3	1.89	0.43
1:D:203[A]:MET:HE1	1:D:217:LEU:HD11	2.01	0.43
1:B:137:ASN:HD21	3:B:402:TXD:H2N1	1.81	0.43
1:A:123:VAL:HG11	1:A:132:LEU:HD21	2.01	0.42
1:A:276:ILE:HG22	1:A:286:VAL:HG23	2.01	0.42
1:B:242:LYS:NZ	7:B:509:HOH:O	2.52	0.42
1:B:28:GLY:HA3	3:B:402:TXD:O5'	2.19	0.42
1:B:280:TYR:HB3	1:B:312:GLU:HG3	2.00	0.42
1:C:221:LYS:O	1:C:222:ASP:C	2.58	0.42
1:C:163:ASN:HA	1:C:271:PRO:HG2	2.01	0.42
1:A:266:LEU:O	1:C:180:HIS:HB2	2.20	0.41
1:A:105:ARG:O	1:A:109:VAL:N	2.48	0.41
1:B:304:LYS:HE3	1:C:6:GLN:O	2.21	0.41
1:B:217:LEU:HD12	1:B:226:TRP:HB2	2.03	0.41
1:B:286:VAL:HG21	1:B:319:ALA:HB1	2.03	0.41
1:C:2:THR:O	1:C:6:GLN:HG3	2.21	0.41
1:C:12:LEU:HD12	1:C:12:LEU:O	2.20	0.40
1:B:11:LEU:HD23	1:C:154[A]:LYS:HG3	2.03	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	327/331 (99%)	315 (96%)	9 (3%)	3 (1%)	17	8
1	B	323/331 (98%)	317 (98%)	6 (2%)	0	100	100
1	C	322/331 (97%)	312 (97%)	7 (2%)	3 (1%)	17	8
1	D	330/331 (100%)	322 (98%)	6 (2%)	2 (1%)	25	15
All	All	1302/1324 (98%)	1266 (97%)	28 (2%)	8 (1%)	25	15

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	105	ARG
1	C	14	GLU
1	A	97	ALA
1	D	97	ALA
1	C	222	ASP
1	C	97	ALA
1	D	16	GLN
1	A	27	VAL

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	288/287 (100%)	281 (98%)	7 (2%)	49	42
1	B	284/287 (99%)	280 (99%)	4 (1%)	67	65
1	C	283/287 (99%)	279 (99%)	4 (1%)	67	65
1	D	290/287 (101%)	282 (97%)	8 (3%)	43	37
All	All	1145/1148 (100%)	1122 (98%)	23 (2%)	57	50

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	14	GLU
1	A	55[A]	ASP
1	A	55[B]	ASP
1	A	105	ARG
1	A	107	ASN
1	A	234	VAL
1	A	236	SER
1	B	55	ASP
1	B	106	LEU
1	B	165	ASP
1	B	236	SER
1	C	15	GLU
1	C	108	LEU
1	C	165	ASP
1	C	220	ASP
1	D	12	LEU
1	D	14	GLU
1	D	16	GLN
1	D	72[A]	ARG
1	D	72[B]	ARG
1	D	165	ASP
1	D	241	ILE
1	D	275	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	B	16	GLN
1	D	330	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

15 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	6EY	B	403	-	28,31,31	1.42	2 (7%)	29,45,45	2.04	6 (20%)
5	SO4	D	803	-	4,4,4	0.30	0	6,6,6	0.20	0
3	TXD	D	801	-	42,48,48	2.28	12 (28%)	49,73,73	2.20	14 (28%)
2	EPE	A	401	-	15,15,15	1.81	1 (6%)	18,20,20	4.84	9 (50%)
2	EPE	C	401	-	15,15,15	1.73	1 (6%)	18,20,20	5.19	9 (50%)
5	SO4	C	404	-	4,4,4	0.34	0	6,6,6	0.31	0
4	6EY	C	403	-	28,31,31	1.36	2 (7%)	29,45,45	1.61	3 (10%)
4	6EY	A	403	-	28,31,31	1.28	2 (7%)	29,45,45	1.62	5 (17%)
3	TXD	B	402	-	42,48,48	2.34	12 (28%)	49,73,73	2.13	11 (22%)
3	TXD	A	402	-	42,48,48	2.15	11 (26%)	49,73,73	2.14	14 (28%)
4	6EY	D	802	-	28,31,31	1.37	2 (7%)	29,45,45	1.80	5 (17%)
5	SO4	A	404	-	4,4,4	0.15	0	6,6,6	0.11	0
6	NAI	C	402	-	42,48,48	2.35	12 (28%)	47,73,73	2.20	8 (17%)
5	SO4	B	404	-	4,4,4	0.22	0	6,6,6	0.11	0
2	EPE	B	401	-	15,15,15	1.66	1 (6%)	18,20,20	5.00	12 (66%)

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	6EY	B	403	-	-	0/10/34/34	0/4/4/4
3	TXD	D	801	-	-	5/26/72/72	0/5/5/5
2	EPE	A	401	-	-	8/9/19/19	0/1/1/1
2	EPE	C	401	-	-	6/9/19/19	0/1/1/1
4	6EY	C	403	-	-	1/10/34/34	0/4/4/4
4	6EY	A	403	-	-	0/10/34/34	0/4/4/4

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	TXD	B	402	-	-	5/26/72/72	0/5/5/5
3	TXD	A	402	-	-	5/26/72/72	0/5/5/5
4	6EY	D	802	-	-	0/10/34/34	0/4/4/4
6	NAI	C	402	-	-	2/25/72/72	0/5/5/5
2	EPE	B	401	-	-	6/9/19/19	0/1/1/1

All (58) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	C	402	NAI	C6N-C5N	9.47	1.50	1.33
3	B	402	TXD	C4N-C5N	7.30	1.50	1.32
3	A	402	TXD	C4N-C5N	7.15	1.50	1.32
3	D	801	TXD	C4N-C5N	7.15	1.50	1.32
2	A	401	EPE	C10-S	-6.39	1.68	1.77
2	C	401	EPE	C10-S	-5.97	1.69	1.77
3	B	402	TXD	C2N-C3N	-5.93	1.47	1.52
2	B	401	EPE	C10-S	-5.74	1.69	1.77
4	B	403	6EY	C17-S18	5.18	1.77	1.70
6	C	402	NAI	C2N-C3N	4.92	1.48	1.34
4	C	403	6EY	C17-S18	4.88	1.77	1.70
4	D	802	6EY	C17-C16	-4.78	1.34	1.37
3	A	402	TXD	C2N-C3N	-4.77	1.48	1.52
3	D	801	TXD	C2'-C3'	-4.58	1.40	1.53
4	A	403	6EY	C17-S18	4.53	1.76	1.70
3	B	402	TXD	C2'-C3'	-4.44	1.41	1.53
3	A	402	TXD	C2'-C3'	-4.42	1.41	1.53
3	D	801	TXD	C2N-C3N	-4.37	1.48	1.52
3	D	801	TXD	C2N-N1N	-4.32	1.40	1.47
3	B	402	TXD	C2N-N1N	-4.31	1.40	1.47
3	D	801	TXD	C7N-N7N	4.24	1.43	1.32
4	B	403	6EY	C17-C16	-4.22	1.34	1.37
3	B	402	TXD	C7N-N7N	4.22	1.43	1.32
6	C	402	NAI	C2B-C3B	-4.15	1.42	1.53
4	C	403	6EY	C17-C16	-3.98	1.34	1.37
3	D	801	TXD	O2'A-C2'A	-3.84	1.33	1.43
3	A	402	TXD	C7N-N7N	3.83	1.42	1.32
6	C	402	NAI	C7N-N7N	3.82	1.43	1.33
4	D	802	6EY	C17-S18	3.62	1.75	1.70
4	A	403	6EY	C17-C16	-3.59	1.34	1.37
3	A	402	TXD	C2N-N1N	-3.51	1.41	1.47
6	C	402	NAI	O2D-C2D	-3.41	1.34	1.43

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	402	TXD	O2'A-C2'A	-3.38	1.35	1.43
3	B	402	TXD	C6A-N6A	3.17	1.45	1.34
6	C	402	NAI	C6A-N6A	3.17	1.45	1.34
3	D	801	TXD	C6A-N6A	2.98	1.44	1.34
3	A	402	TXD	C6A-N6A	2.95	1.44	1.34
6	C	402	NAI	C6N-N1N	2.88	1.44	1.37
6	C	402	NAI	C5D-C4D	-2.84	1.42	1.51
3	A	402	TXD	C5'A-C4'A	-2.82	1.42	1.51
3	D	801	TXD	C3N-C7N	-2.74	1.49	1.53
3	A	402	TXD	O2'A-C2'A	-2.73	1.36	1.43
3	D	801	TXD	C5'A-C4'A	-2.72	1.43	1.51
3	B	402	TXD	C3N-C7N	-2.50	1.50	1.53
3	D	801	TXD	C3'A-C2'A	-2.40	1.46	1.53
3	B	402	TXD	C2'-C1'	-2.40	1.50	1.53
3	B	402	TXD	C5'A-C4'A	-2.38	1.44	1.51
3	D	801	TXD	C2'-C1'	-2.32	1.50	1.53
6	C	402	NAI	C2B-C1B	-2.29	1.50	1.53
6	C	402	NAI	O3D-C3D	-2.28	1.37	1.43
6	C	402	NAI	C2D-C3D	-2.14	1.47	1.53
3	A	402	TXD	C2'-C1'	-2.09	1.50	1.53
3	B	402	TXD	C5'-C4'	-2.06	1.45	1.51
6	C	402	NAI	O4B-C4B	-2.06	1.40	1.45
3	A	402	TXD	C3N-C7N	-2.06	1.50	1.53
3	D	801	TXD	O2'-C2'	-2.04	1.38	1.43
3	A	402	TXD	C3'A-C2'A	-2.04	1.47	1.53
3	B	402	TXD	O4'-C4'	-2.01	1.40	1.45

All (96) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	401	EPE	O1S-S-C10	-14.21	89.81	106.92
2	C	401	EPE	O2S-S-C10	-14.02	90.04	106.92
2	A	401	EPE	O1S-S-C10	-12.60	91.75	106.92
2	C	401	EPE	O3S-S-C10	-10.45	88.87	105.77
2	C	401	EPE	O1S-S-C10	-9.78	95.13	106.92
2	A	401	EPE	O2S-S-C10	-9.38	95.62	106.92
6	C	402	NAI	C1D-N1N-C2N	8.52	135.30	121.11
2	A	401	EPE	O3S-S-C10	-8.28	92.38	105.77
2	B	401	EPE	O2S-S-C10	-7.85	97.46	106.92
4	B	403	6EY	C15-N14-C13	7.25	122.73	117.33
3	B	402	TXD	O4'A-C1'A-N1N	7.20	119.83	109.47
6	C	402	NAI	C1D-N1N-C6N	-7.17	105.38	120.83

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	801	TXD	C2N-C3N-C4N	6.64	118.11	108.63
3	A	402	TXD	C2N-C3N-C4N	6.45	117.84	108.63
3	D	801	TXD	O4'A-C1'A-N1N	6.42	118.72	109.47
3	B	402	TXD	C3N-C4N-C5N	-6.28	112.54	123.09
2	B	401	EPE	O3S-S-C10	-6.00	96.07	105.77
4	C	403	6EY	C15-N14-C13	5.91	121.73	117.33
3	D	801	TXD	C3N-C4N-C5N	-5.87	113.22	123.09
3	A	402	TXD	C3N-C4N-C5N	-5.82	113.31	123.09
3	A	402	TXD	O4'A-C1'A-N1N	5.38	117.21	109.47
4	D	802	6EY	C16-C17-S18	-5.37	108.83	112.29
6	C	402	NAI	O4D-C1D-N1N	5.28	118.37	108.06
3	B	402	TXD	C4N-C3N-C7N	5.12	124.25	110.08
2	A	401	EPE	C6-N1-C2	5.08	120.26	108.83
2	C	401	EPE	C6-N1-C2	4.98	120.04	108.83
4	A	403	6EY	C15-N14-C13	4.96	121.02	117.33
4	B	403	6EY	C16-C17-S18	-4.93	109.11	112.29
3	B	402	TXD	C2N-C3N-C4N	4.83	115.52	108.63
2	A	401	EPE	C6-C5-N4	-4.75	100.90	110.64
2	B	401	EPE	C6-N1-C2	4.73	119.47	108.83
3	A	402	TXD	N3A-C2A-N1A	-4.67	121.38	128.68
2	B	401	EPE	C6-C5-N4	4.64	120.16	110.64
2	B	401	EPE	C5-C6-N1	4.59	120.06	110.64
6	C	402	NAI	N3A-C2A-N1A	-4.45	121.73	128.68
2	B	401	EPE	C2-C3-N4	4.44	119.74	110.64
2	A	401	EPE	O3S-S-O1S	4.40	122.02	111.27
2	B	401	EPE	C5-N4-C3	4.37	118.66	108.83
3	A	402	TXD	C4N-C3N-C7N	4.22	121.77	110.08
4	D	802	6EY	C15-N14-C13	4.17	120.43	117.33
3	B	402	TXD	N3A-C2A-N1A	-4.08	122.29	128.68
2	B	401	EPE	C3-C2-N1	4.01	118.86	110.64
3	D	801	TXD	N3A-C2A-N1A	-3.97	122.47	128.68
4	A	403	6EY	C16-C17-S18	-3.96	109.74	112.29
3	D	801	TXD	C4N-C3N-C7N	3.81	120.64	110.08
2	C	401	EPE	O3S-S-O2S	3.73	120.38	111.27
3	B	402	TXD	C5N-C6N-N1N	3.55	119.49	111.03
4	C	403	6EY	C16-C17-S18	-3.53	110.01	112.29
3	D	801	TXD	C5N-C6N-N1N	3.50	119.38	111.03
4	D	802	6EY	C4-C13-N14	3.49	120.09	114.39
4	C	403	6EY	C4-C13-N14	3.48	120.08	114.39
3	D	801	TXD	O4'-C1'-C2'	-3.46	101.87	106.93
2	A	401	EPE	C7-N4-C5	3.42	119.98	111.23
2	A	401	EPE	O3S-S-O2S	3.41	119.61	111.27

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	802	6EY	C6-S5-C4	3.40	107.13	102.81
4	A	403	6EY	C3-C15-N14	3.30	112.08	108.09
2	C	401	EPE	O3S-S-O1S	3.25	119.22	111.27
2	A	401	EPE	C3-C2-N1	3.21	117.23	110.64
2	C	401	EPE	C5-C6-N1	3.15	117.11	110.64
3	D	801	TXD	PA-O3-PN	-3.06	122.33	132.83
3	A	402	TXD	C6N-N1N-C1'A	-2.99	104.08	112.83
6	C	402	NAI	C4D-O4D-C1D	-2.95	102.97	109.47
4	B	403	6EY	C4-C13-N14	2.93	119.19	114.39
3	A	402	TXD	C5N-C6N-N1N	2.91	117.96	111.03
3	A	402	TXD	C4'A-O4'A-C1'A	-2.91	103.05	109.47
3	B	402	TXD	PA-O3-PN	-2.88	122.94	132.83
2	C	401	EPE	C2-C3-N4	-2.84	104.81	110.64
2	B	401	EPE	O3S-S-O1S	2.78	118.07	111.27
3	D	801	TXD	C4'A-O4'A-C1'A	-2.78	103.35	109.47
6	C	402	NAI	PN-O3-PA	-2.77	123.32	132.83
3	D	801	TXD	C4A-C5A-N7A	-2.75	106.54	109.40
3	D	801	TXD	C6N-N1N-C1'A	-2.70	104.91	112.83
3	B	402	TXD	O4'-C1'-C2'	-2.65	103.06	106.93
3	A	402	TXD	C3N-C7N-N7N	2.65	119.70	116.85
4	A	403	6EY	C4-C13-N14	2.64	118.71	114.39
2	C	401	EPE	C7-N4-C3	2.61	117.90	111.23
2	B	401	EPE	O2S-S-O1S	2.60	122.94	113.95
3	A	402	TXD	O4'-C1'-C2'	-2.53	103.24	106.93
4	B	403	6EY	C24-C25-C27	-2.50	117.44	120.17
3	A	402	TXD	PA-O3-PN	-2.50	124.26	132.83
3	D	801	TXD	C3N-C7N-N7N	2.48	119.52	116.85
4	D	802	6EY	C3-C15-N14	2.44	111.04	108.09
2	B	401	EPE	O3S-S-O2S	2.44	117.23	111.27
3	A	402	TXD	O7N-C7N-N7N	-2.36	118.90	123.00
3	B	402	TXD	O5'A-C5'A-C4'A	2.35	117.08	108.99
3	D	801	TXD	O5'-C5'-C4'	2.28	116.84	108.99
4	B	403	6EY	C3-C15-N14	2.24	110.80	108.09
3	B	402	TXD	C6N-N1N-C1'A	-2.24	106.26	112.83
3	B	402	TXD	C4'A-O4'A-C1'A	-2.23	104.56	109.47
4	B	403	6EY	C27-C21-C15	-2.17	117.25	120.74
6	C	402	NAI	C4A-C5A-N7A	-2.12	107.19	109.40
4	A	403	6EY	C6-S5-C4	2.11	105.48	102.81
3	A	402	TXD	C3'A-C2'A-C1'A	-2.07	97.50	101.43
6	C	402	NAI	O4B-C1B-C2B	-2.04	103.95	106.93
3	A	402	TXD	C4A-C5A-N7A	-2.02	107.30	109.40
3	D	801	TXD	O5'A-C5'A-C4'A	2.01	115.91	108.99

There are no chirality outliers.

All (38) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	D	801	TXD	C2N-C3N-C7N-O7N
3	D	801	TXD	C2N-C3N-C7N-N7N
3	D	801	TXD	O4'A-C1'A-N1N-C2N
2	A	401	EPE	C8-C7-N4-C5
2	A	401	EPE	C9-C10-S-O1S
2	A	401	EPE	C9-C10-S-O3S
2	C	401	EPE	C8-C7-N4-C3
3	B	402	TXD	C2N-C3N-C7N-N7N
3	B	402	TXD	O4'A-C1'A-N1N-C2N
3	A	402	TXD	C2N-C3N-C7N-O7N
3	A	402	TXD	C2N-C3N-C7N-N7N
3	A	402	TXD	O4'A-C1'A-N1N-C2N
2	B	401	EPE	C10-C9-N1-C6
2	C	401	EPE	C9-C10-S-O3S
2	B	401	EPE	C9-C10-S-O3S
3	B	402	TXD	C2N-C3N-C7N-O7N
2	A	401	EPE	S-C10-C9-N1
2	A	401	EPE	C10-C9-N1-C2
2	A	401	EPE	C10-C9-N1-C6
2	C	401	EPE	C8-C7-N4-C5
2	B	401	EPE	N4-C7-C8-O8
2	A	401	EPE	C9-C10-S-O2S
2	C	401	EPE	C9-C10-S-O1S
2	C	401	EPE	C9-C10-S-O2S
2	B	401	EPE	C9-C10-S-O1S
2	B	401	EPE	C9-C10-S-O2S
3	D	801	TXD	C2'A-C1'A-N1N-C2N
3	A	402	TXD	C2'A-C1'A-N1N-C2N
2	A	401	EPE	C8-C7-N4-C3
2	C	401	EPE	C10-C9-N1-C6
2	B	401	EPE	C10-C9-N1-C2
6	C	402	NAI	O4D-C1D-N1N-C2N
3	B	402	TXD	C4N-C3N-C7N-N7N
3	D	801	TXD	O4'-C4'-C5'-O5'
3	B	402	TXD	O4'-C4'-C5'-O5'
6	C	402	NAI	O4B-C4B-C5B-O5B
3	A	402	TXD	O4'-C4'-C5'-O5'
4	C	403	6EY	C16-C15-C21-C27

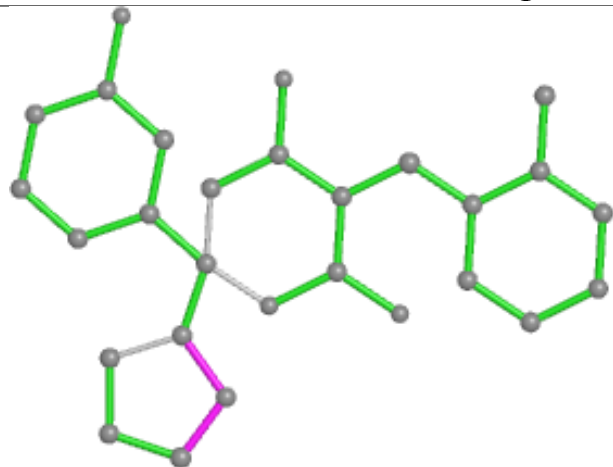
There are no ring outliers.

7 monomers are involved in 20 short contacts:

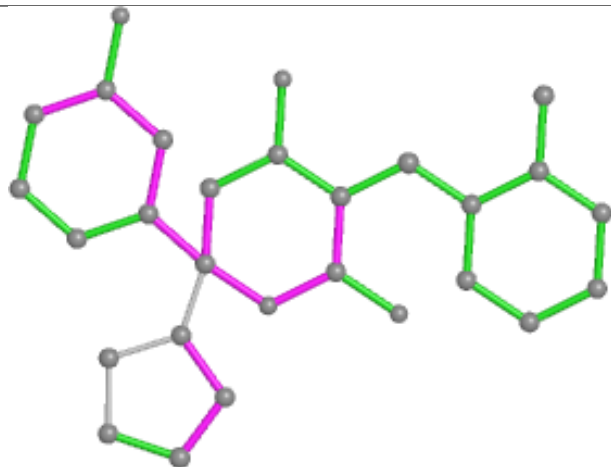
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	403	6EY	2	0
3	D	801	TXD	2	0
4	A	403	6EY	2	0
3	B	402	TXD	6	0
3	A	402	TXD	6	0
6	C	402	NAI	1	0
2	B	401	EPE	1	0

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.

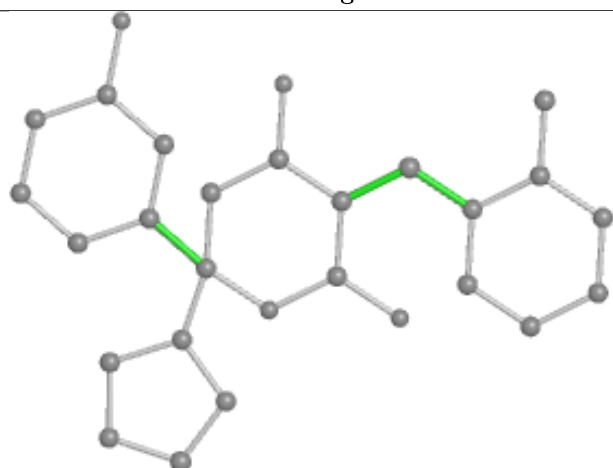
Ligand 6EY B 403



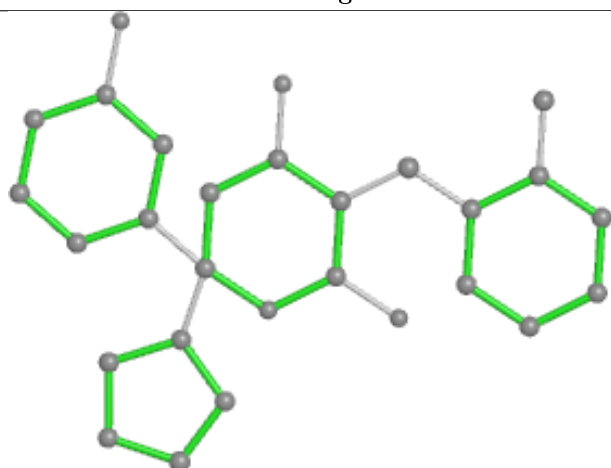
Bond lengths



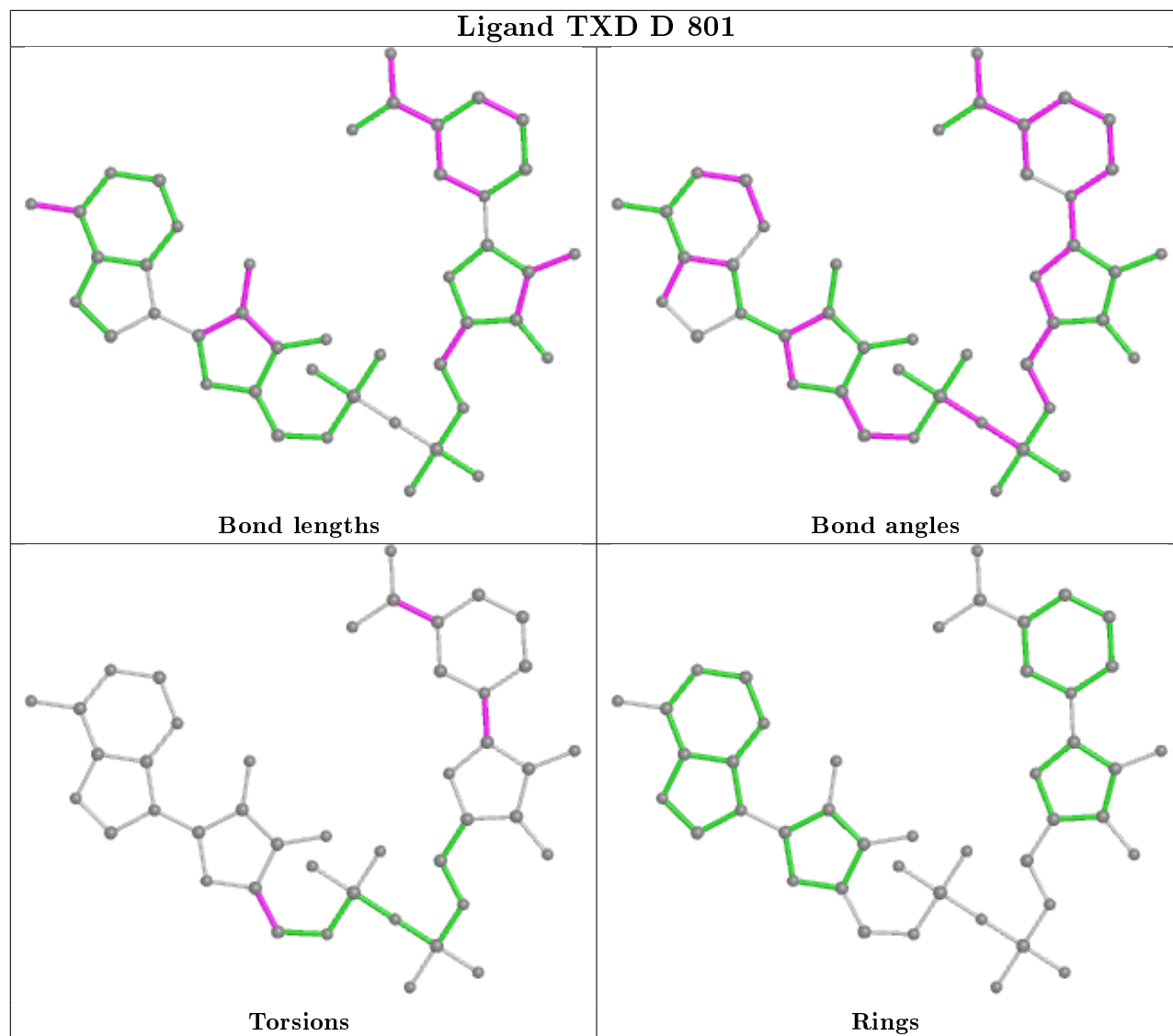
Bond angles



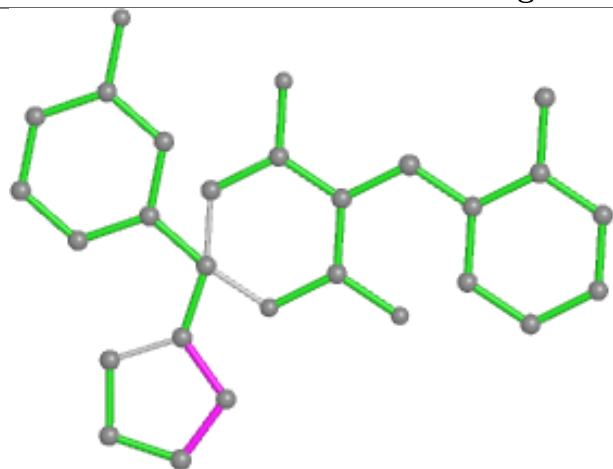
Torsions



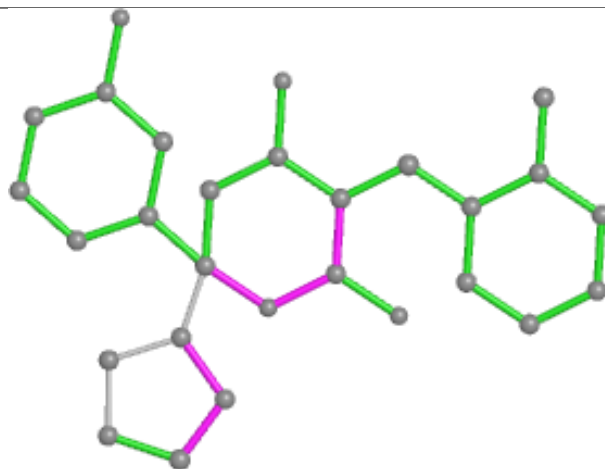
Rings



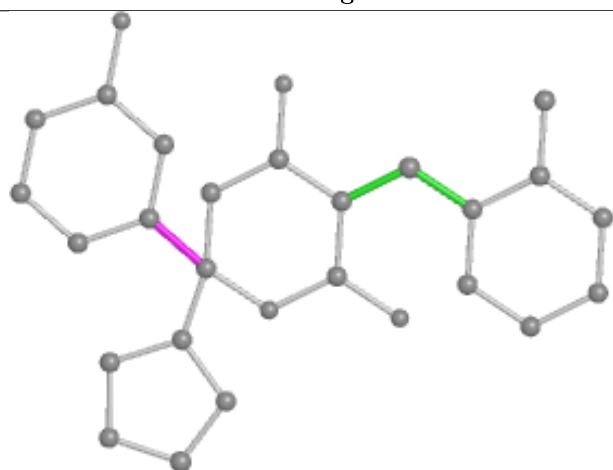
Ligand 6EY C 403



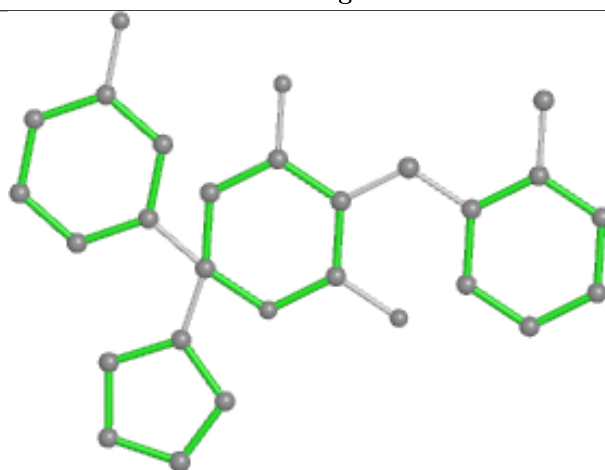
Bond lengths



Bond angles

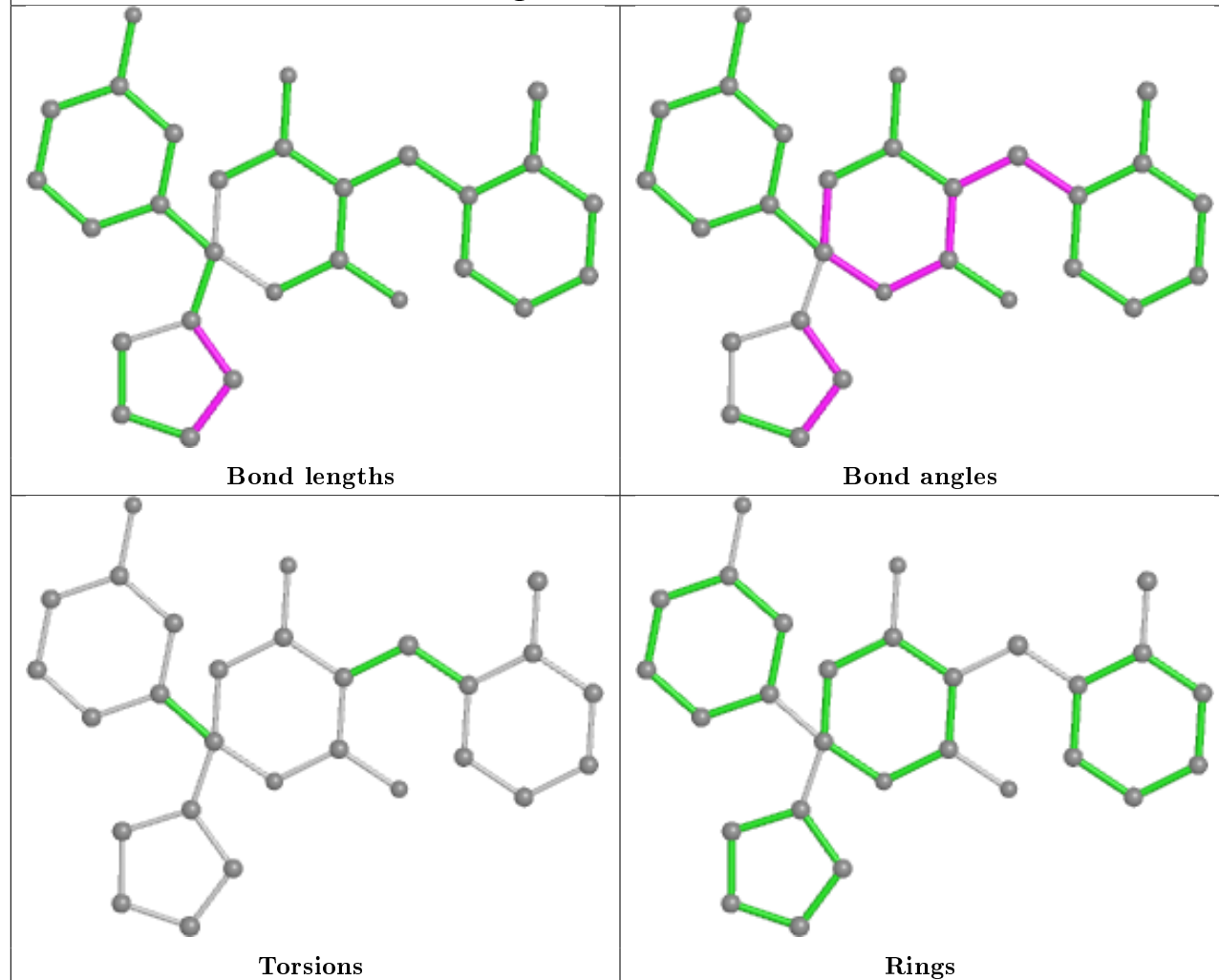


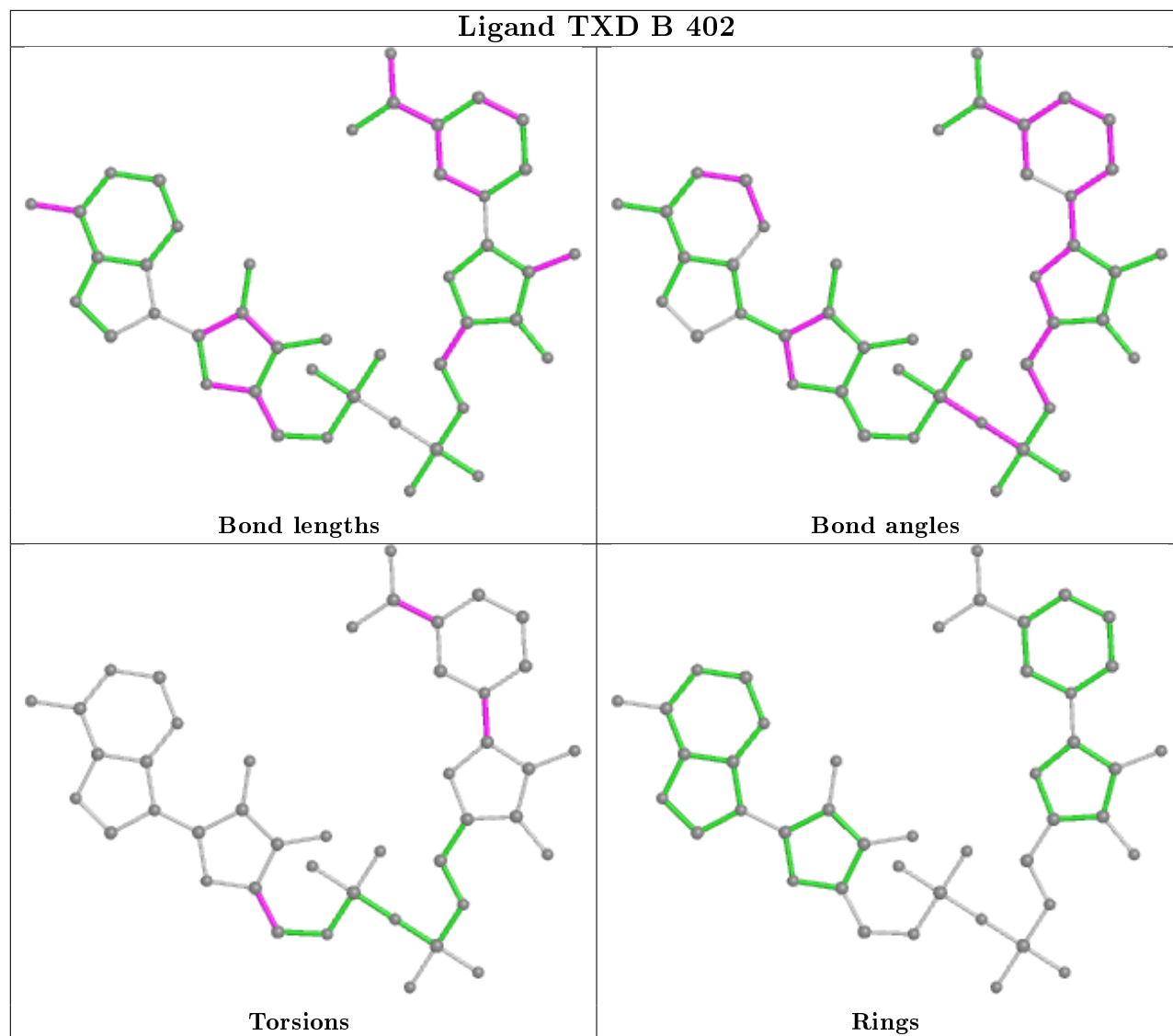
Torsions

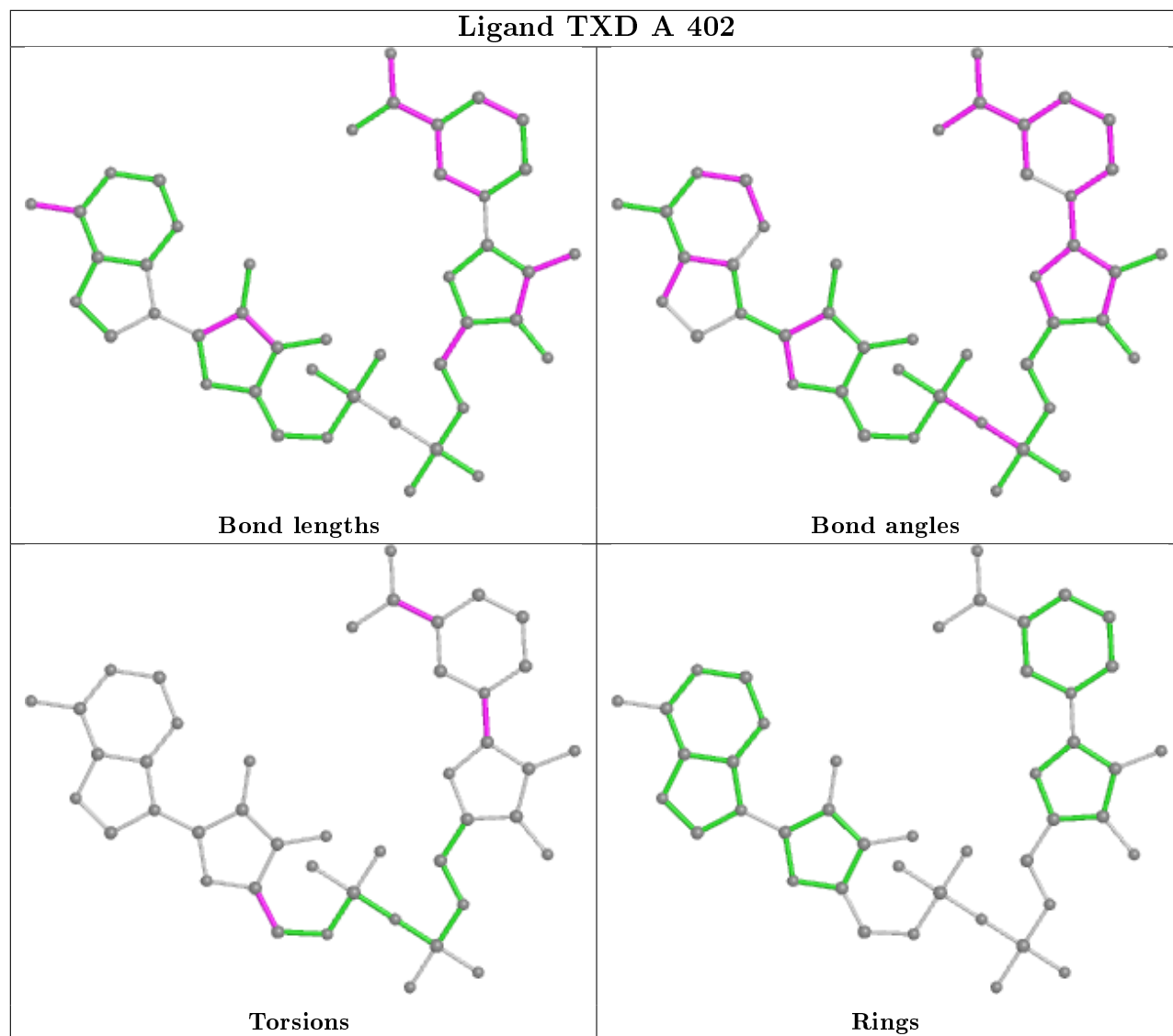


Rings

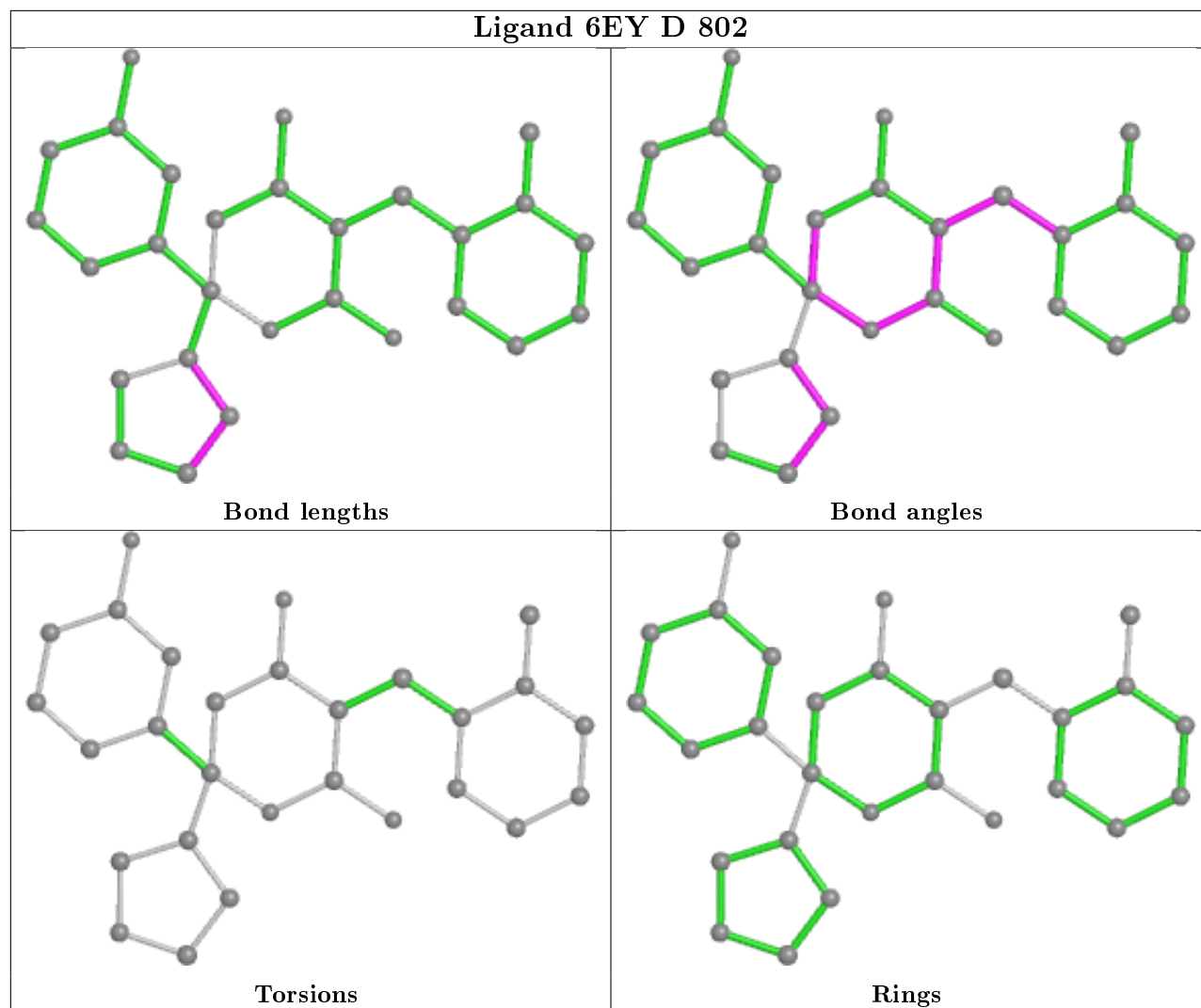
Ligand 6EY A 403

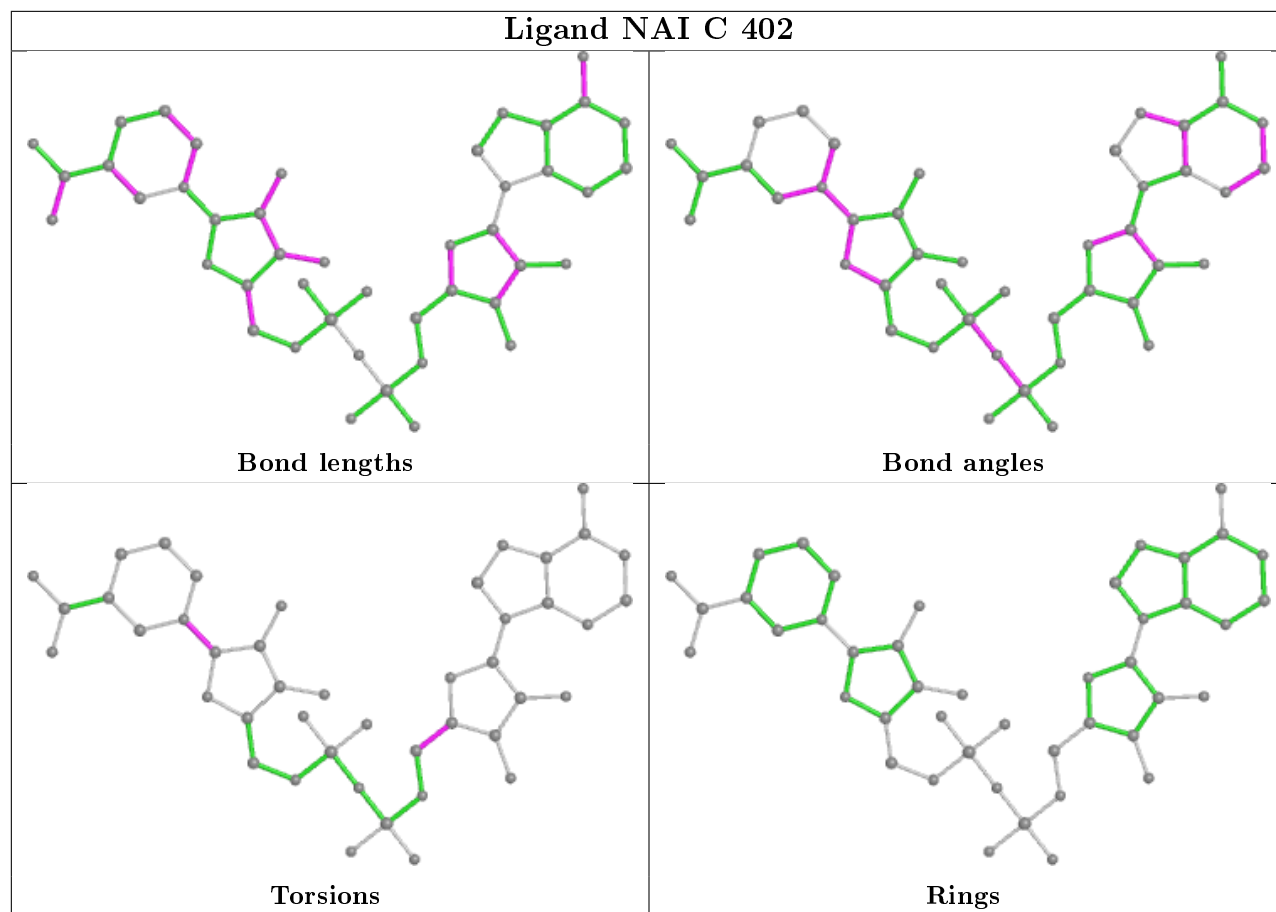






Ligand 6EY D 802





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	327/331 (98%)	-0.28	2 (0%) 89 91	18, 30, 59, 88	0
1	B	324/331 (97%)	-0.27	8 (2%) 57 61	19, 30, 60, 108	0
1	C	324/331 (97%)	0.04	19 (5%) 22 24	21, 38, 83, 135	0
1	D	330/331 (99%)	-0.32	10 (3%) 50 54	20, 30, 68, 119	0
All	All	1305/1324 (98%)	-0.21	39 (2%) 50 54	18, 32, 71, 135	0

All (39) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	97	ALA	7.4
1	B	108	LEU	4.9
1	C	14	GLU	4.7
1	D	15	GLU	4.3
1	D	1	ALA	4.1
1	C	329	LEU	3.7
1	D	100	GLN	3.5
1	D	107	ASN	3.2
1	C	107	ASN	3.1
1	D	104	SER	3.1
1	C	330	GLN	3.1
1	D	331	PHE	3.1
1	A	97	ALA	3.0
1	C	110	GLN	2.9
1	C	2	THR	2.9
1	D	13	LYS	2.9
1	B	13	LYS	2.8
1	A	105	ARG	2.8
1	C	13	LYS	2.8
1	D	106	LEU	2.8
1	B	106	LEU	2.7

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	C	327	LYS	2.6
1	C	108	LEU	2.6
1	C	331	PHE	2.5
1	B	330	GLN	2.5
1	B	72[A]	ARG	2.5
1	B	107	ASN	2.5
1	C	328	GLU	2.3
1	C	221	LYS	2.3
1	B	331	PHE	2.3
1	C	106	LEU	2.3
1	C	128	PRO	2.3
1	C	111	ARG	2.2
1	C	98	ARG	2.2
1	B	324	GLY	2.2
1	D	16	GLN	2.1
1	D	72[A]	ARG	2.0
1	C	114	ASN	2.0
1	C	220	ASP	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 carbohydrates in this entry.

6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	EPE	C	401	15/15	0.77	0.21	49,76,92,95	15
2	EPE	A	401	15/15	0.84	0.19	34,57,65,67	15
4	6EY	B	403	28/28	0.91	0.17	34,57,72,83	0
6	NAI	C	402	44/44	0.92	0.14	34,50,60,68	0
2	EPE	B	401	15/15	0.93	0.15	43,57,81,94	15

Continued on next page...

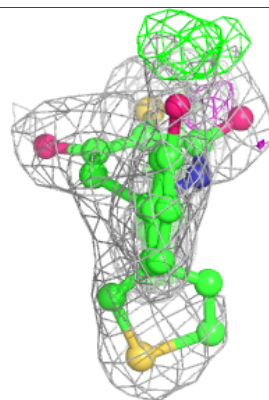
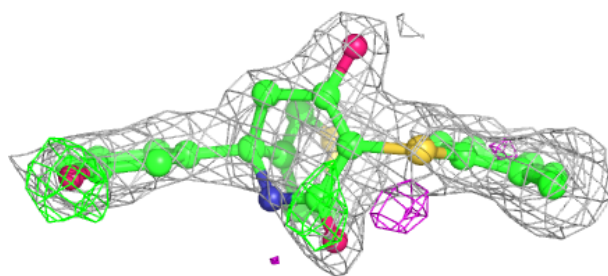
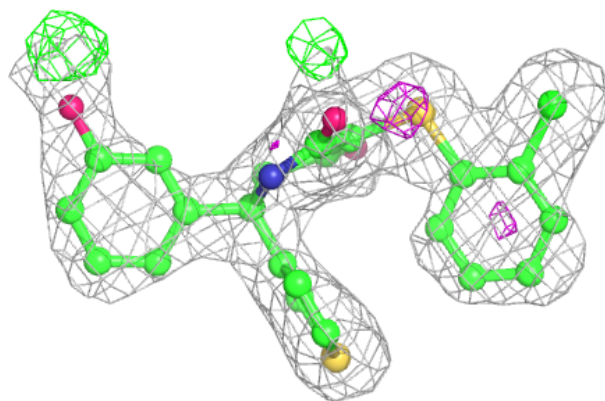
Continued from previous page...

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
4	6EY	C	403	28/28	0.94	0.13	26,51,71,84	0
4	6EY	A	403	28/28	0.94	0.11	26,37,47,57	0
4	6EY	D	802	28/28	0.95	0.12	31,38,59,71	0
3	TXD	B	402	44/44	0.96	0.10	24,34,40,45	0
3	TXD	D	801	44/44	0.96	0.09	22,34,40,53	0
3	TXD	A	402	44/44	0.97	0.08	21,30,38,44	0
5	SO4	A	404	5/5	0.99	0.18	23,24,29,35	4
5	SO4	C	404	5/5	0.99	0.15	24,24,31,36	4
5	SO4	B	404	5/5	0.99	0.17	20,24,26,32	4
5	SO4	D	803	5/5	0.99	0.16	21,23,26,32	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.

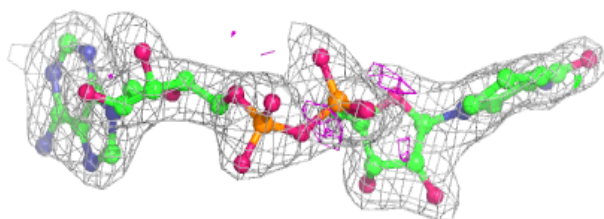
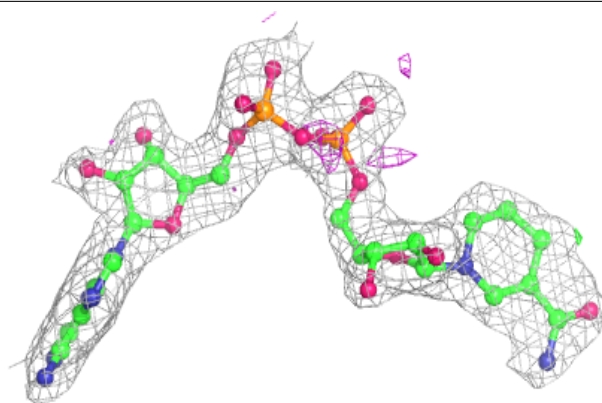
Electron density around 6EY B 403:

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

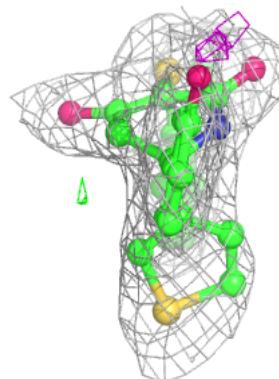
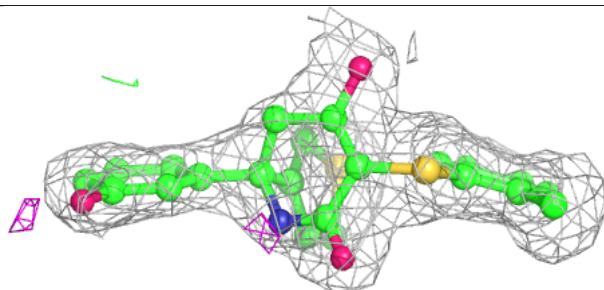
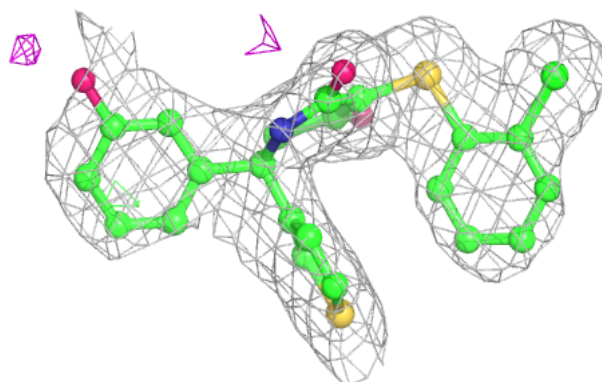


Electron density around NAI C 402:

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

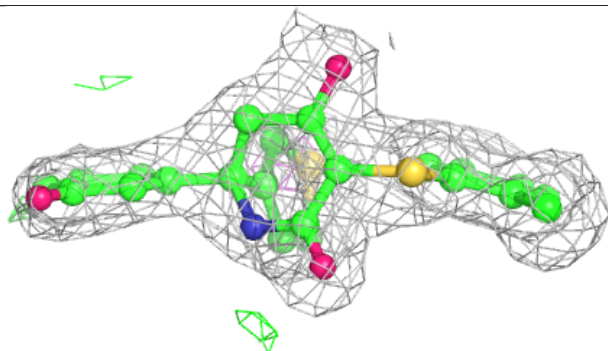
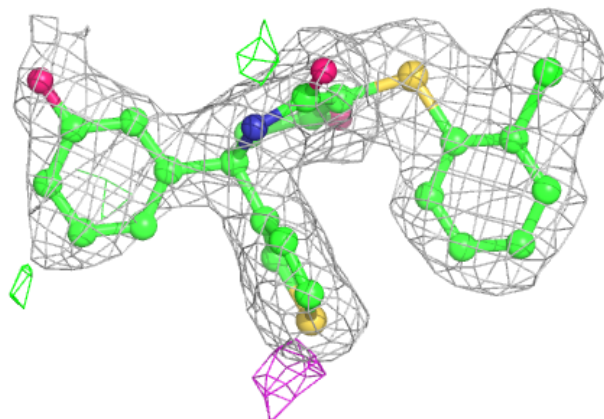
**Electron density around 6EY C 403:**

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

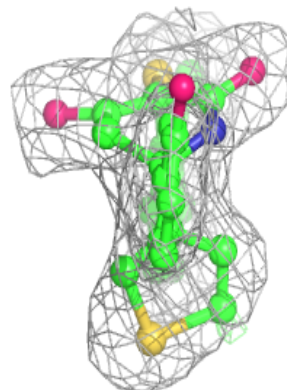
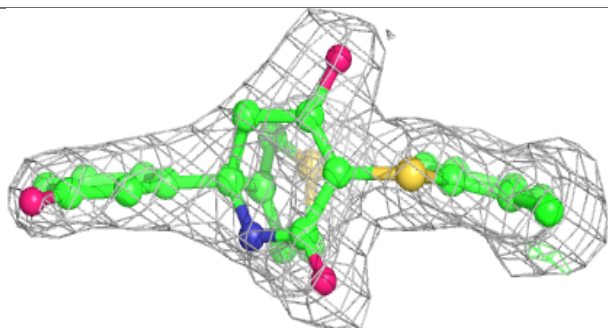
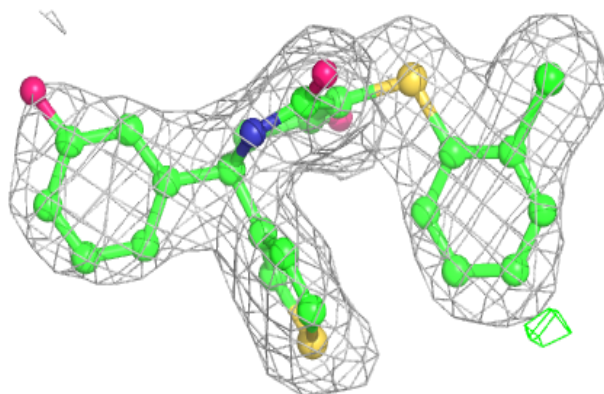


Electron density around 6EY A 403:

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

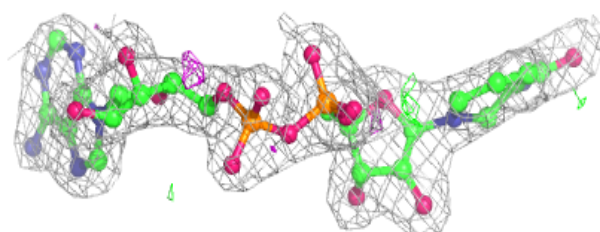
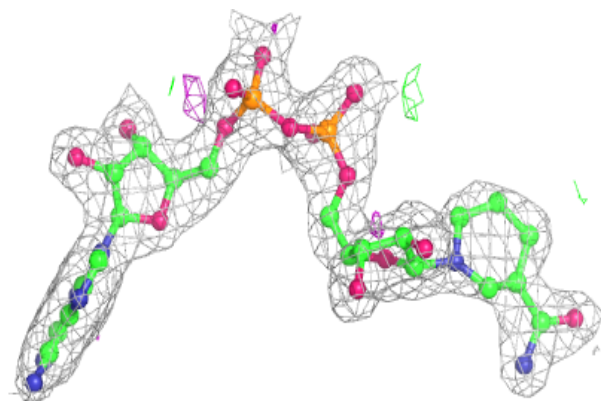
**Electron density around 6EY D 802:**

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

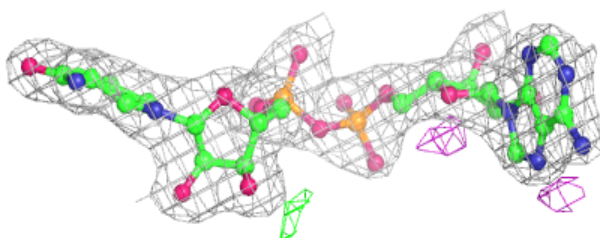
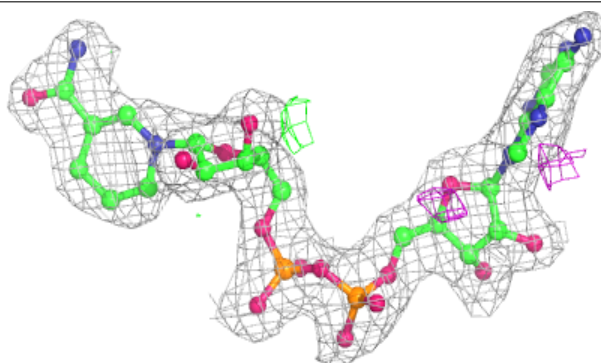


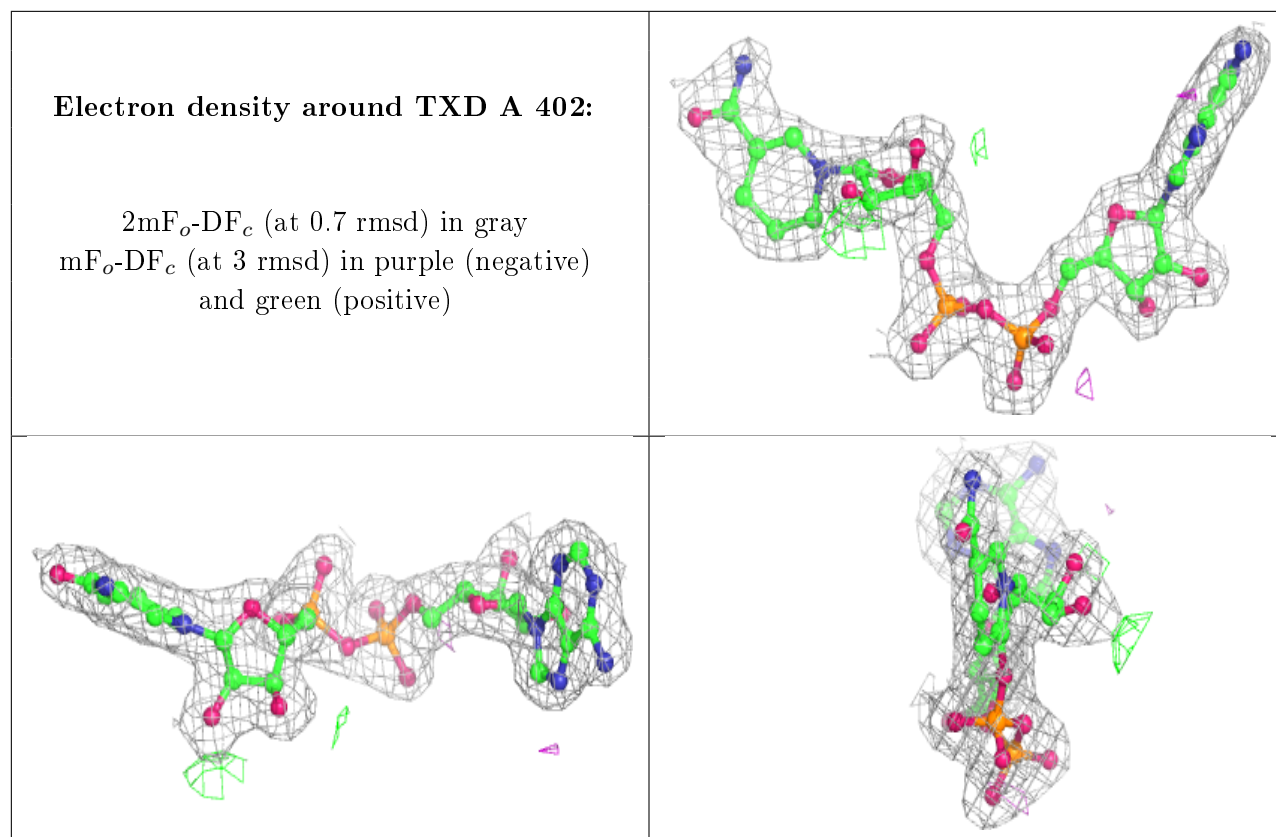
Electron density around TXD B 402:

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

**Electron density around TXD D 801:**

$2mF_o-DF_c$ (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.