

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP070720\
 Data File : BP002662.D
 Acq On : 07 Jul 2020 17:31
 Operator : CG/JU
 Sample : L3205-01
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 12013

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0
 Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP062920.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.193	385	392	409	rBV	3385480	5227127	59.22%	10.233%
2	6.763	652	659	674	rBV	3353653	5569456	63.10%	10.903%
3	7.569	789	796	806	rBV	824811	1286957	14.58%	2.519%
4	7.887	841	850	861	rBV	3193659	5237872	59.34%	10.254%
5	8.716	974	991	1004	rBV	2338520	4039565	45.77%	7.908%
6	10.340	1260	1267	1273	rBV	1020517	1742791	19.75%	3.412%
7	10.387	1273	1275	1283	rVB	57446	90346	1.02%	0.177%
8	12.016	1547	1552	1562	rBV2	65885	154987	1.76%	0.303%
9	12.234	1585	1589	1595	rVB	65787	100127	1.13%	0.196%
10	12.645	1654	1659	1665	rBV6	51179	115298	1.31%	0.226%
11	12.834	1684	1691	1707	rBV	5531329	8826459	100.00%	17.279%
12	13.004	1716	1720	1724	rBV3	68916	111662	1.27%	0.219%
13	13.134	1737	1742	1749	rBV	340023	570267	6.46%	1.116%
14	13.645	1824	1829	1832	rBV	356711	564014	6.39%	1.104%
15	13.681	1832	1835	1841	rVB	453276	626468	7.10%	1.226%
16	13.887	1867	1870	1874	rBV3	119050	163491	1.85%	0.320%
17	13.963	1880	1883	1888	rBV4	158317	217227	2.46%	0.425%
18	14.057	1895	1899	1904	rVB	489360	673568	7.63%	1.319%
19	14.216	1919	1926	1934	rBV	1420233	2650352	30.03%	5.188%
20	15.022	2060	2063	2070	rBV3	491064	717487	8.13%	1.405%
21	15.722	2178	2182	2189	rBV	2797930	4219380	47.80%	8.260%
22	19.574	2834	2837	2842	rVB	3000918	3949354	44.74%	7.731%
23	21.092	3092	3095	3099	rVB	1421967	1698703	19.25%	3.325%
24	23.286	3463	3468	3474	rVB	1490081	2529008	28.65%	4.951%

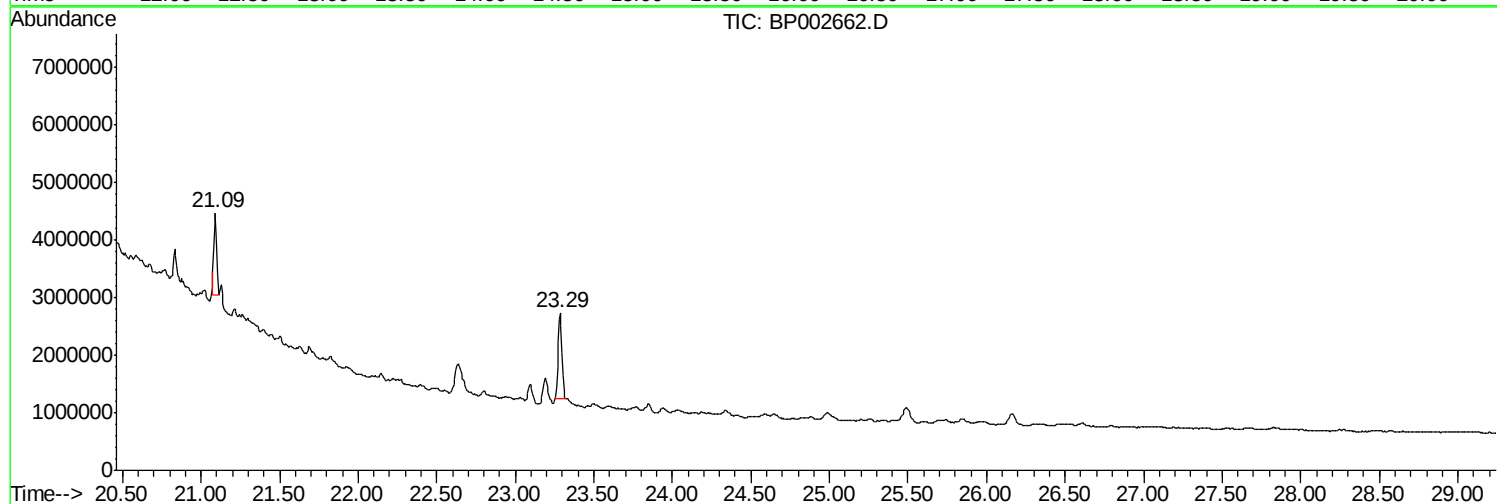
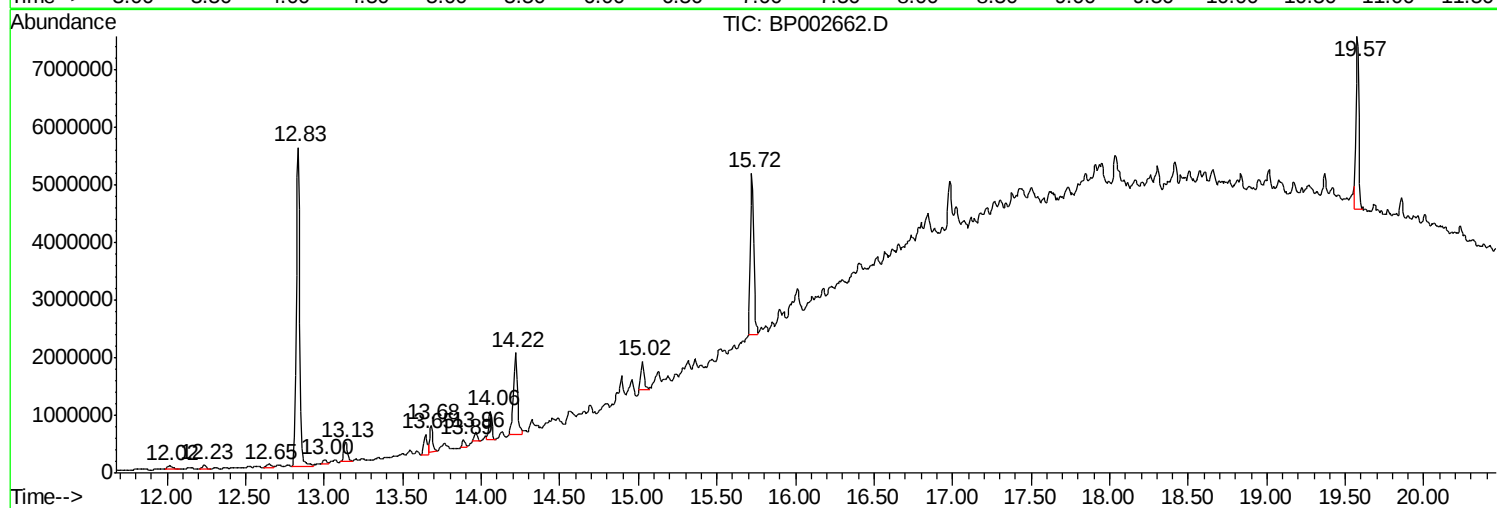
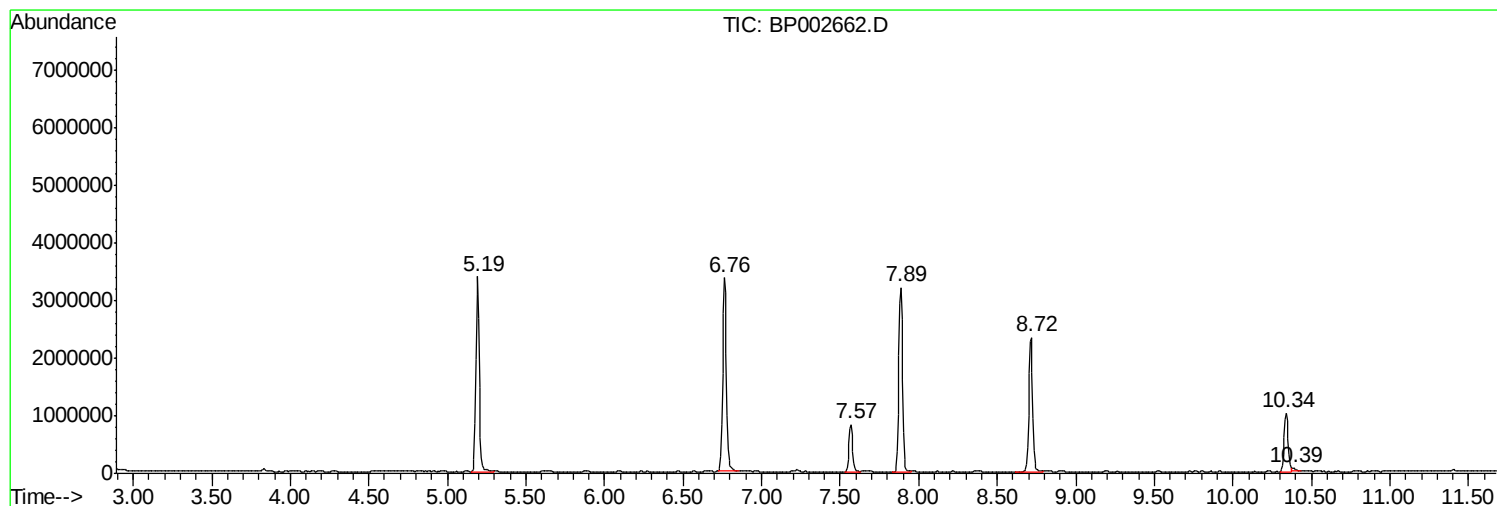
Sum of corrected areas: 51081966

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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP062920.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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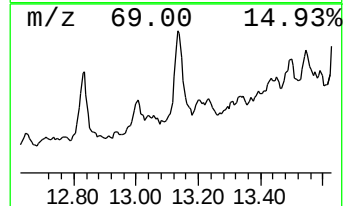
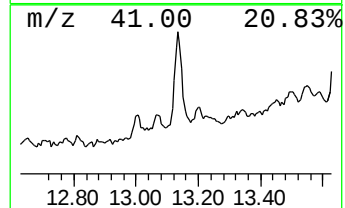
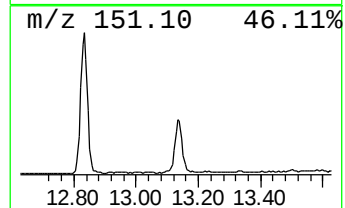
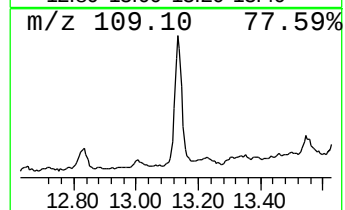
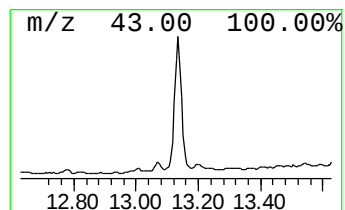
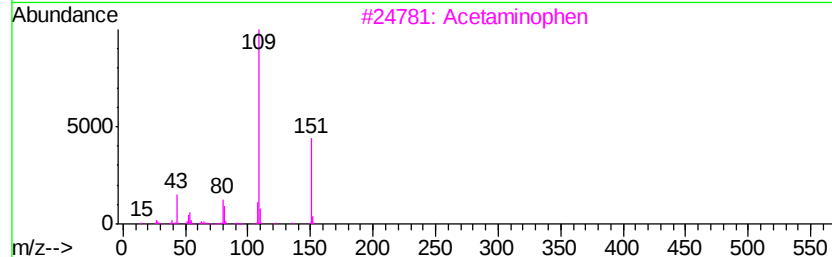
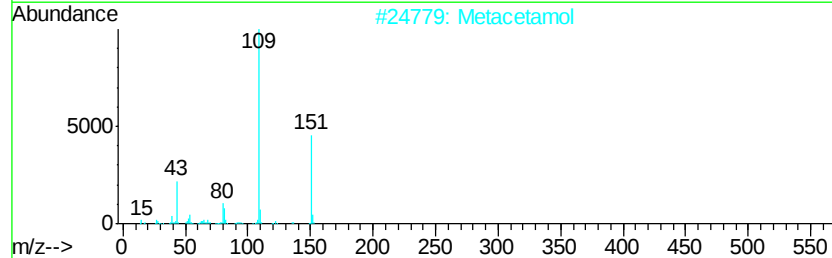
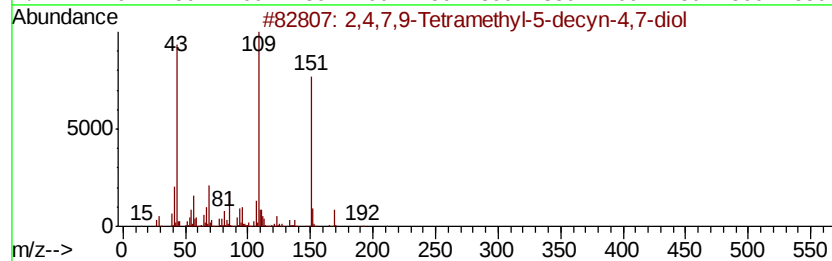
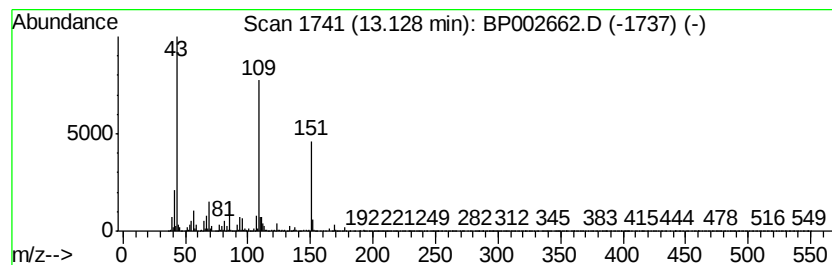
Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP062920.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 2,4,7,9-Tetramethyl-5-decyn... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.13	4.30 ng	570267	Acenaphthene-d10	14.22

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4,7,9-Tetramethyl-5-decyn-4,7-...	226	C14H26O2	000126-86-3	91		
2	Metacetamol	151	C8H9NO2	000621-42-1	64		
3	Acetaminophen	151	C8H9NO2	000103-90-2	59		
4	Pvridine, 2-butyl-, 1-oxide	151	C9H13NO	031396-32-4	59		
5	m-Isopropoxyaniline	151	C9H13NO	041406-00-2	59		



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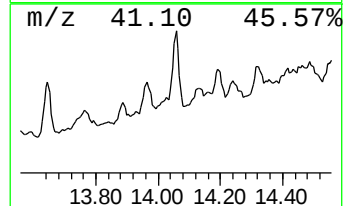
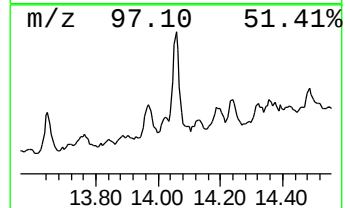
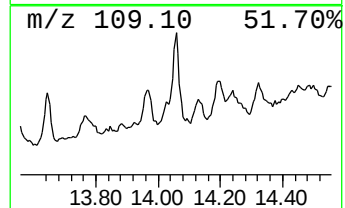
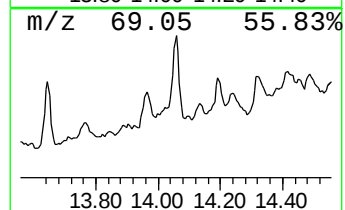
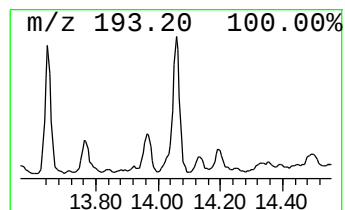
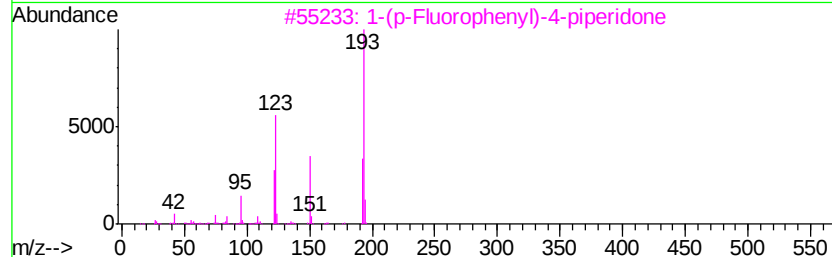
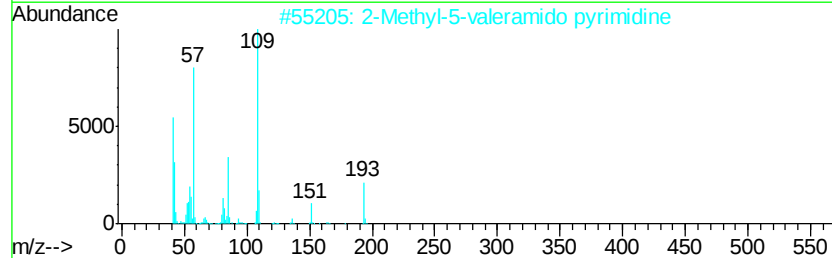
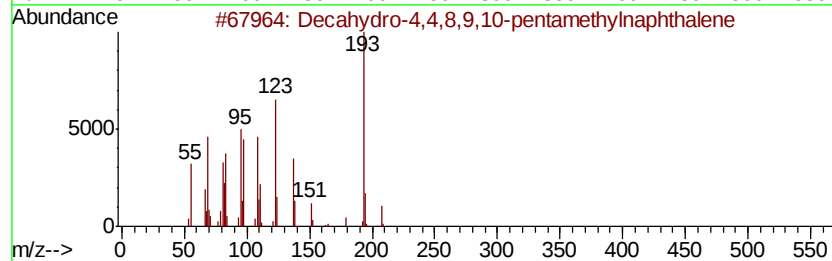
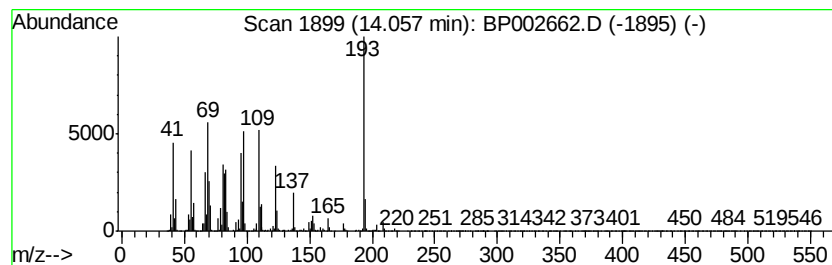
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Peak Number 3 unknown14.06 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.06	5.08 ng	673568	Acenaphthene-d10	14.22
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Decahydro-4,4,8,9,10-pentamethyl...	208 C15H28	080655-44-3 49
2		2-Methyl-5-valeramido pyrimidine	193 C10H15N3O	1000213-95-9 35
3		1-(p-Fluorophenyl)-4-piperidone	193 C11H12FN0	1000238-56-7 32
4		Silane, chlorodiethyl(2-methylpe...	222 C10H23Cl0Si	1000363-79-1 25
5		Borinic acid, diethyl-, 3,3,5-tr...	208 C13H25B0	057387-76-5 25



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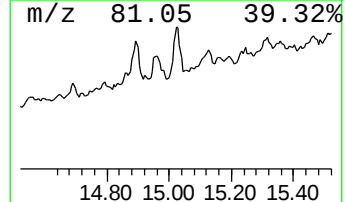
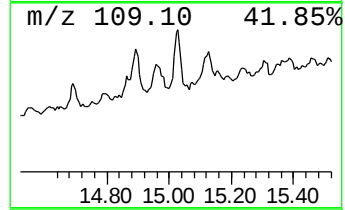
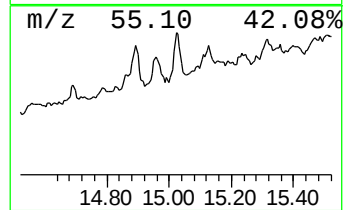
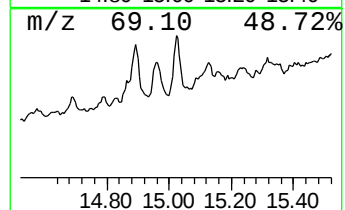
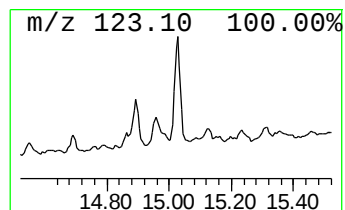
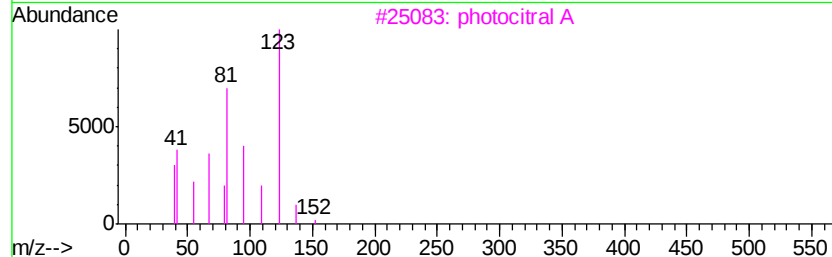
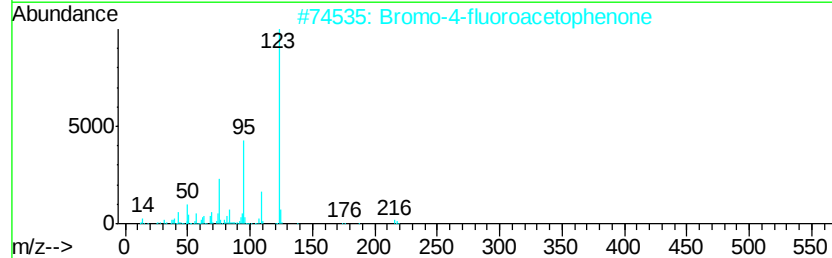
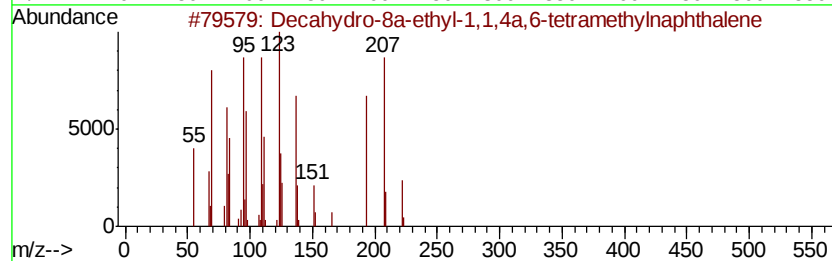
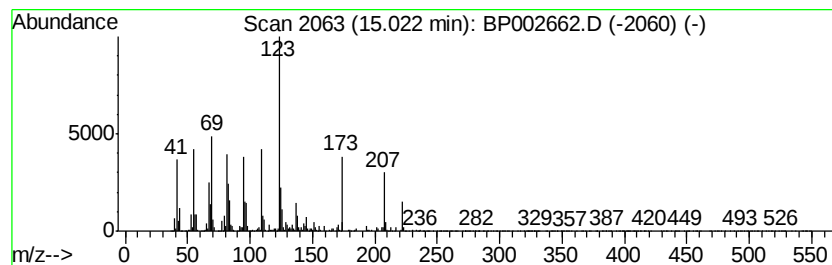
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Peak Number 4 unknown15.02 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.02	5.41 ng	717487	Acenaphthene-d10	14.22

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Decahydro-8a-ethyl-1,1,4a,6-tetr...	222	C16H30	1000100-23-6	43
2		Bromo-4-fluoroacetophenone	216	C8H6BrFO	000403-29-2	38
3		photocitral A	152	C10H16O	055253-28-6	38
4		3-Fluorobenzoic acid, 3-methylbu...	208	C12H13FO2	154559-38-3	35
5		2,6-Naphthalenedione, octahydro-...	208	C13H20O2	057289-17-5	32



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2,4,7,9-Tetrameth...	13.13	4.3	ng	570267	3	14.22	2650350	20.0
unknown14.06	14.06	5.1	ng	673568	3	14.22	2650350	20.0
unknown15.02	15.02	5.4	ng	717487	3	14.22	2650350	20.0