

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM022721\
 Data File : BM028944.D
 Acq On : 27 Feb 2021 06:53
 Operator : CG/JU
 Sample : M1498-06
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 CB8F9

Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : OFF
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 0.1 % 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_M\METHODS\SOM-EPA-BM022721MA.M
 Title : SVOA CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.122	20	23	27	rBV	4640	5463	1.14%	0.100%
2	3.157	27	29	34	rVB	2171	2444	0.51%	0.045%
3	4.351	228	232	235	rBB	2829	3368	0.70%	0.062%
4	5.045	345	350	356	rBB	30606	45317	9.43%	0.829%
5	6.957	670	675	684	rVB	26353	42433	8.83%	0.776%
6	7.110	696	701	709	rBB	78847	117608	24.49%	2.150%
7	7.304	728	734	741	rBV	97572	149727	31.17%	2.737%
8	7.410	747	752	758	rBB3	1174	2414	0.50%	0.044%
9	7.769	808	813	824	rBB	63472	99856	20.79%	1.826%
10	8.139	871	876	880	rBB2	2764	4103	0.85%	0.075%
11	8.504	932	938	952	rBB	73732	116881	24.33%	2.137%
12	8.945	1007	1013	1023	rBB	105491	170946	35.59%	3.125%
13	9.286	1068	1071	1074	rBV	1649	2115	0.44%	0.039%
14	9.398	1086	1090	1094	rBB	4227	5543	1.15%	0.101%
15	9.663	1130	1135	1143	rBB	95974	153391	31.94%	2.804%
16	10.098	1204	1209	1215	rBB2	1143	2051	0.43%	0.037%
17	10.204	1221	1227	1238	rBB	137767	220966	46.00%	4.040%
18	10.569	1283	1289	1296	rBV	85590	140715	29.30%	2.573%
19	10.727	1309	1316	1324	rBV2	9034	15512	3.23%	0.284%
20	11.927	1512	1520	1526	rBB2	5764	10320	2.15%	0.189%
21	12.098	1543	1549	1552	rBV2	1211	1848	0.38%	0.034%
22	12.133	1552	1555	1558	rBV2	1629	2284	0.48%	0.042%
23	12.169	1558	1561	1564	rVB2	2049	2259	0.47%	0.041%
24	12.369	1591	1595	1598	rBV3	1511	2163	0.45%	0.040%
25	12.463	1604	1611	1617	rVB5	2688	4623	0.96%	0.085%
26	12.533	1617	1623	1628	rVB3	1913	4107	0.86%	0.075%
27	12.586	1628	1632	1636	rBV5	1392	2407	0.50%	0.044%
28	12.633	1636	1640	1644	rVB4	1397	2126	0.44%	0.039%
29	12.951	1688	1694	1701	rVB4	1735	3949	0.82%	0.072%
30	13.145	1723	1727	1729	rBV4	1641	2411	0.50%	0.044%
31	13.239	1737	1743	1747	rBV3	1883	3743	0.78%	0.068%
32	13.280	1747	1750	1752	rVV2	1959	2499	0.52%	0.046%
33	13.321	1752	1757	1762	rVV	14392	21139	4.40%	0.386%
34	13.392	1764	1769	1773	rBV5	1546	2371	0.49%	0.043%

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35	13.492	1782	1786	1789	rVB3	1693	2660	0.55%	0.049%
36	13.545	1789	1795	1798	rBV4	1240	3035	0.63%	0.055%
37	13.668	1812	1816	1820	rVB4	3793	5192	1.08%	0.095%
38	13.757	1826	1831	1834	rBV3	2159	3262	0.68%	0.060%
39	13.851	1841	1847	1852	rVV	233918	312620	65.09%	5.716%
40	13.898	1852	1855	1862	rVB	8718	11247	2.34%	0.206%
41	13.986	1867	1870	1874	rBV3	2242	4001	0.83%	0.073%
42	14.080	1882	1886	1889	rBV3	5922	7717	1.61%	0.141%
43	14.127	1889	1894	1903	rVB	286445	397918	82.84%	7.275%
44	14.263	1912	1917	1924	rBV	7162	13377	2.79%	0.245%
45	14.363	1931	1934	1939	rVB4	2361	3278	0.68%	0.060%
46	14.433	1940	1946	1952	rBV	128796	185628	38.65%	3.394%
47	14.498	1952	1957	1964	rVB	28620	39346	8.19%	0.719%
48	14.586	1968	1972	1979	rVB	2716	4484	0.93%	0.082%
49	14.680	1984	1988	1998	rBV	19420	34211	7.12%	0.625%
50	14.833	2009	2014	2019	rBV	9108	11832	2.46%	0.216%
51	14.915	2024	2028	2030	rBV3	2508	3394	0.71%	0.062%
52	14.986	2038	2040	2046	rVB3	3129	4659	0.97%	0.085%
53	15.051	2046	2051	2056	rBV6	2703	6683	1.39%	0.122%
54	15.186	2070	2074	2083	rVV2	7330	16797	3.50%	0.307%
55	15.257	2084	2086	2093	rVB3	3387	4847	1.01%	0.089%
56	15.339	2097	2100	2102	rVV3	1626	2368	0.49%	0.043%
57	15.433	2110	2116	2121	rVV	324758	465394	96.89%	8.509%
58	15.480	2121	2124	2130	rVB2	13160	19541	4.07%	0.357%
59	15.574	2130	2140	2149	rBV	123919	175190	36.47%	3.203%
60	15.662	2151	2155	2158	rVV3	7772	9412	1.96%	0.172%
61	15.698	2158	2161	2167	rVB5	3142	5275	1.10%	0.096%
62	15.892	2190	2194	2197	rBV4	1245	1938	0.40%	0.035%
63	15.957	2197	2205	2210	rBV	54012	75523	15.72%	1.381%
64	16.104	2226	2230	2233	rBV5	2375	3553	0.74%	0.065%
65	16.145	2233	2237	2245	rVV3	16236	29788	6.20%	0.545%
66	16.198	2245	2246	2253	rVB6	2691	3247	0.68%	0.059%
67	16.280	2259	2260	2264	rBV3	1882	2355	0.49%	0.043%
68	16.368	2271	2275	2276	rBV2	2466	2649	0.55%	0.048%
69	16.392	2276	2279	2283	rVB2	2589	3239	0.67%	0.059%
70	16.468	2285	2292	2297	rBV	28652	41023	8.54%	0.750%
71	16.751	2339	2340	2347	rVB6	2092	2403	0.50%	0.044%

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72	16.956	2372	2375	2378	rBV4	1625	2507	0.52%	0.046%
73	16.992	2379	2381	2384	rVB2	3789	3323	0.69%	0.061%
74	17.027	2386	2387	2390	rBV3	1491	1839	0.38%	0.034%
75	17.086	2395	2397	2401	rBV5	1510	1967	0.41%	0.036%
76	17.180	2404	2413	2418	rBV	143146	201260	41.90%	3.680%
77	17.221	2418	2420	2423	rVV2	2308	2335	0.49%	0.043%
78	17.280	2423	2430	2438	rVB	328342	469927	97.84%	8.592%
79	17.462	2457	2461	2462	rBV4	2163	2271	0.47%	0.042%
80	17.556	2475	2477	2481	rVB5	2057	2546	0.53%	0.047%
81	17.662	2493	2495	2498	rVB4	3926	3351	0.70%	0.061%
82	17.786	2513	2516	2521	rVV7	1924	3392	0.71%	0.062%
83	17.945	2539	2543	2546	rVB5	1834	2608	0.54%	0.048%
84	18.021	2554	2556	2559	rVB4	2260	1987	0.41%	0.036%
85	18.068	2559	2564	2572	rBV	39551	53654	11.17%	0.981%
86	18.245	2591	2594	2596	rVV3	2431	2606	0.54%	0.048%
87	18.551	2641	2646	2651	rVB4	4764	8110	1.69%	0.148%
88	18.809	2685	2690	2695	rBV	32727	41790	8.70%	0.764%
89	18.898	2701	2705	2709	rBV2	8882	10981	2.29%	0.201%
90	19.203	2753	2757	2760	rBV2	4107	6364	1.32%	0.116%
91	19.345	2777	2781	2788	rBV2	19312	25533	5.32%	0.467%
92	19.445	2795	2798	2803	rBV2	10226	12105	2.52%	0.221%
93	19.562	2812	2818	2824	rBV	349826	480317	100.00%	8.782%
94	19.615	2824	2827	2834	rVB	43305	53513	11.14%	0.978%
95	20.580	2989	2991	2995	rBV4	5419	5894	1.23%	0.108%
96	20.662	3002	3005	3008	rBV	6820	8489	1.77%	0.155%
97	21.044	3067	3070	3079	rVB2	32121	39535	8.23%	0.723%
98	21.339	3115	3120	3125	rBV	140598	173699	36.16%	3.176%
99	23.397	3464	3470	3479	rVB	229572	399175	83.11%	7.298%
100	23.538	3488	3494	3503	rVB	87637	160308	33.38%	2.931%

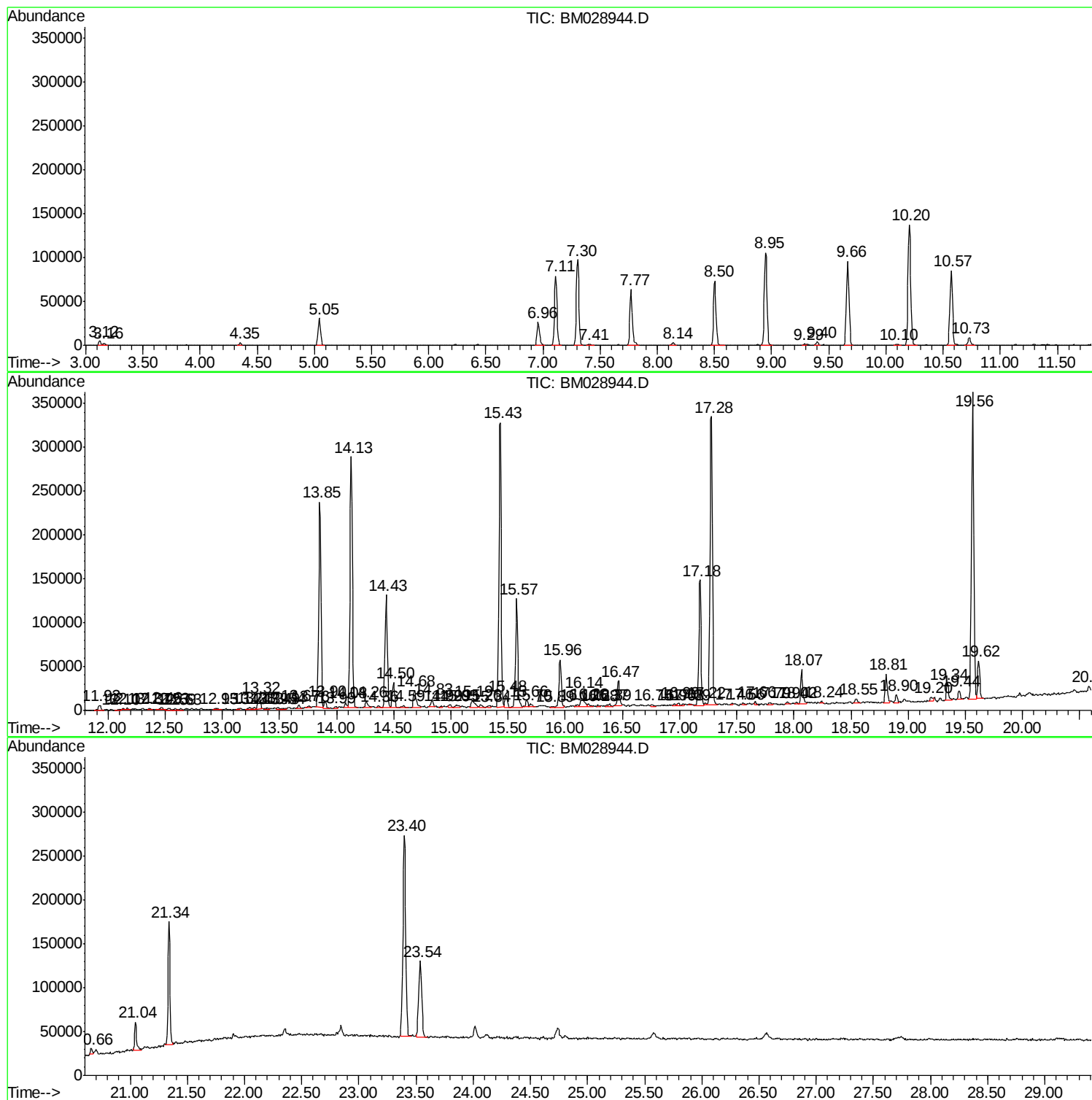
Sum of corrected areas: 5469574

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM022721MA.M
Quant Title : SVOA CALIBRATION

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



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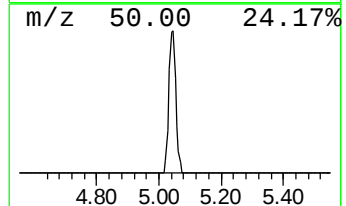
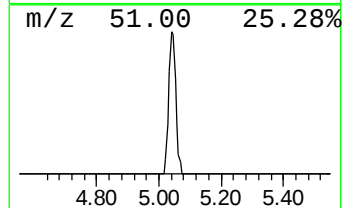
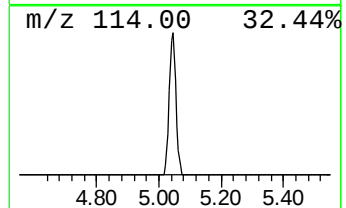
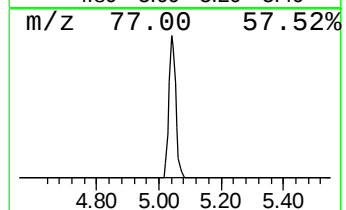
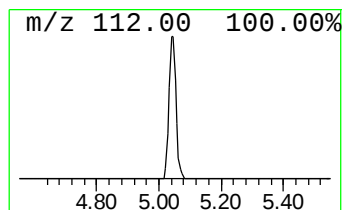
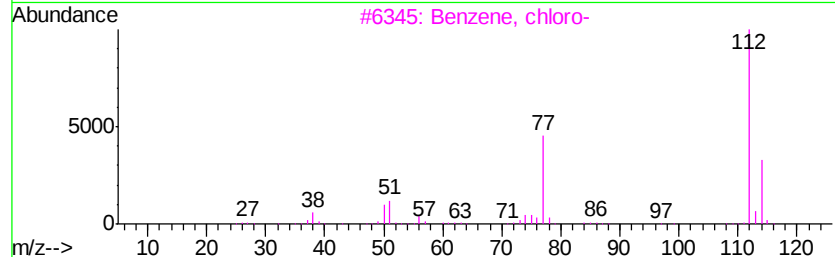
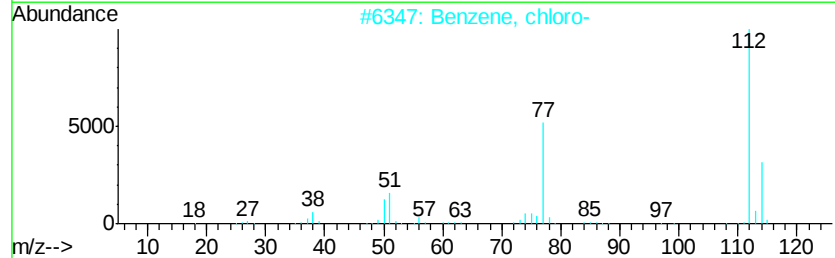
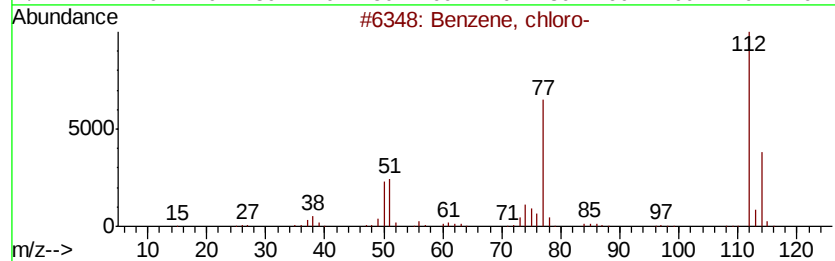
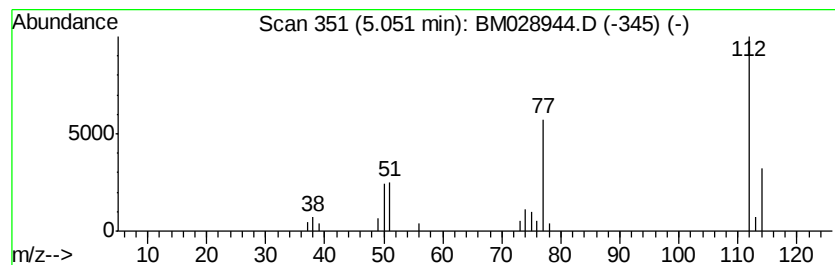
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM022721MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzene, chloro- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.05	9.08 ng/ul	45317	1,4-Dichlorobenzene-d4	7.77

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, chloro-	112	C6H5Cl	000108-90-7	96
2			Benzene, chloro-	112	C6H5Cl	000108-90-7	91
3			Benzene, chloro-	112	C6H5Cl	000108-90-7	91
4			Benzene, chloro-	112	C6H5Cl	000108-90-7	87
5			2-Furancarboxylic acid	112	C5H4O3	000088-14-2	27



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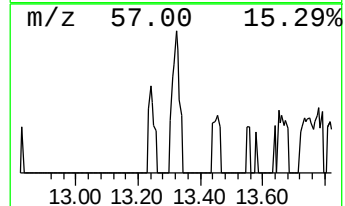
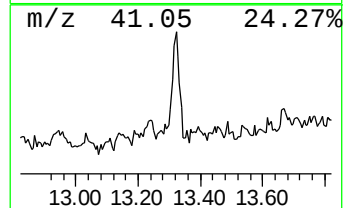
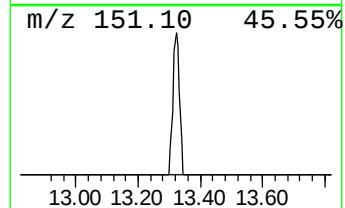
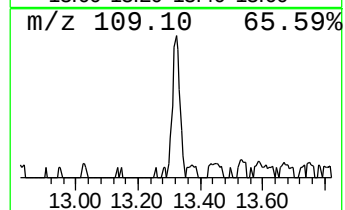
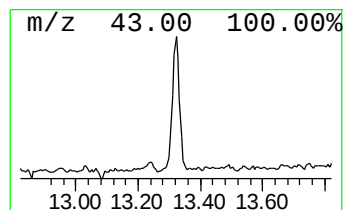
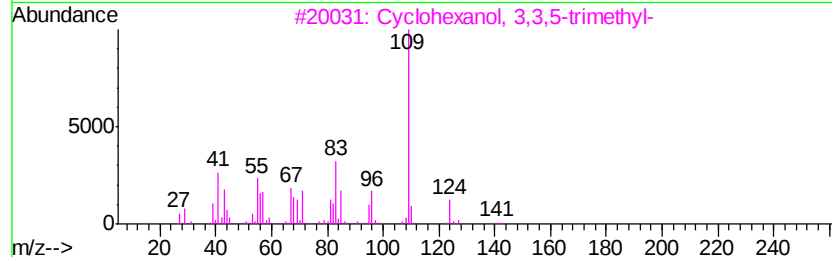
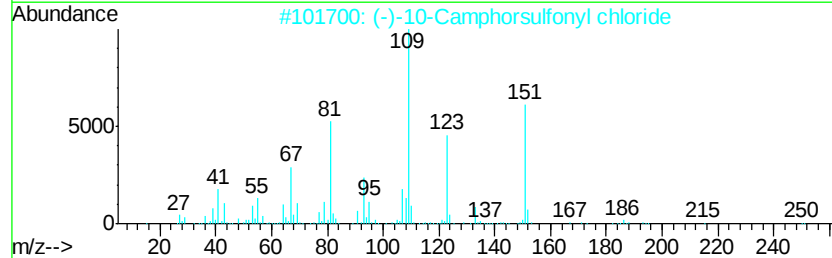
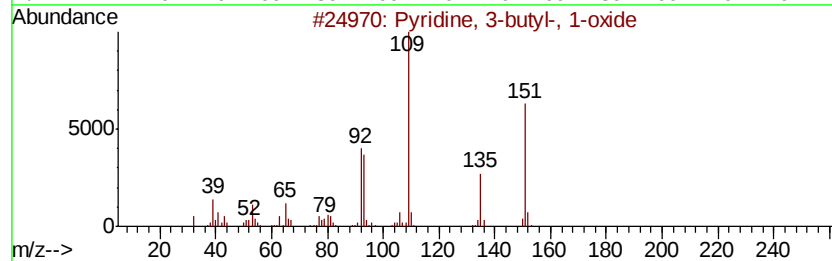
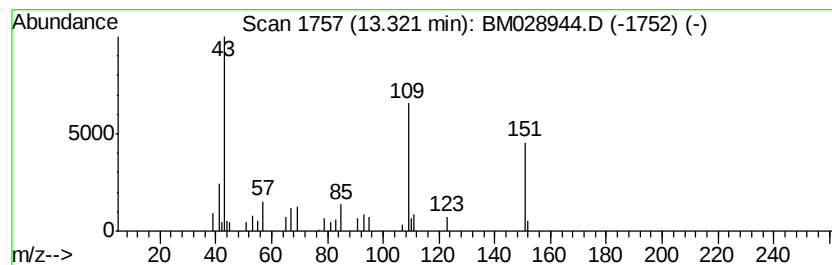
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 2 unknown-01 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.32	2.28 ng/ul	21139	Acenaphthene-d10	14.43
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Pyridine, 3-butyl-, 1-oxide	151 C9H13NO	031396-33-5 42
2		(-)-10-Camphorsulfonyl chloride	250 C10H15ClO3S	039262-22-1 39
3		Cyclohexanol, 3,3,5-trimethyl-	142 C9H18O	000116-02-9 38
4		Ethanone, 1-(1,4-dimethyl-3-cyclo...	152 C10H16O	043219-68-7 38
5		1-Acetyl-3-amino-4-cyano-3-pyrro...	151 C7H9N3O	002125-74-8 32



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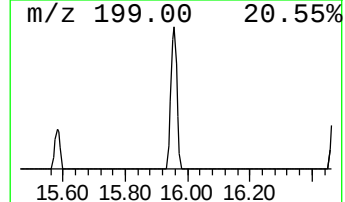
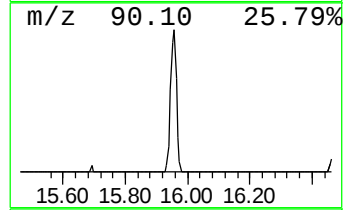
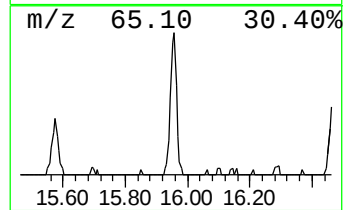
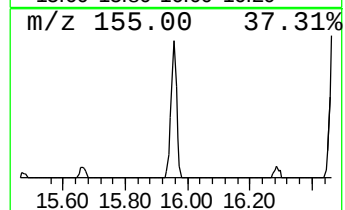
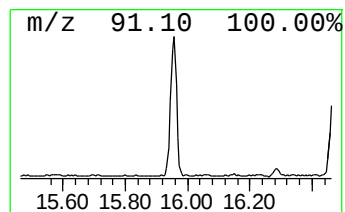
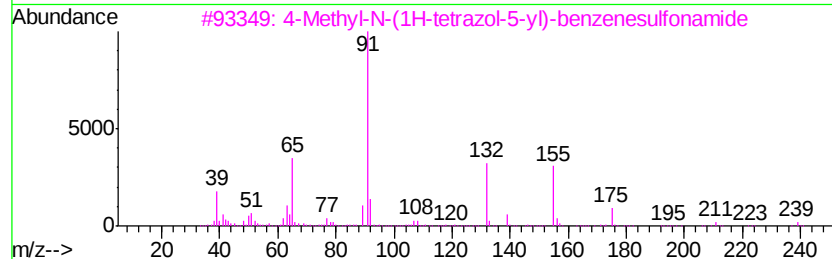
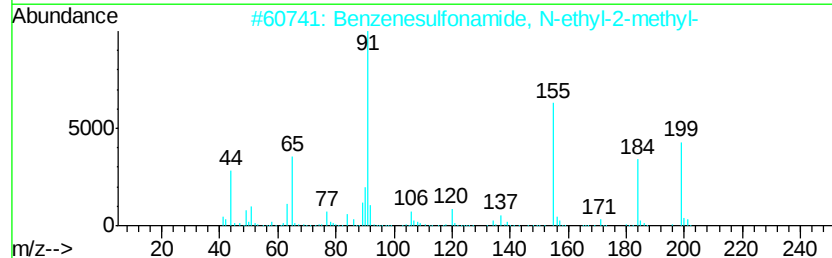
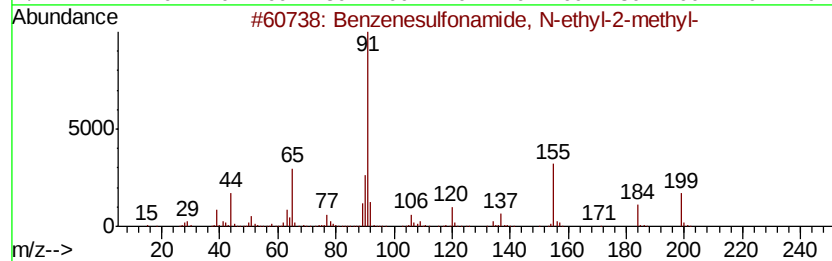
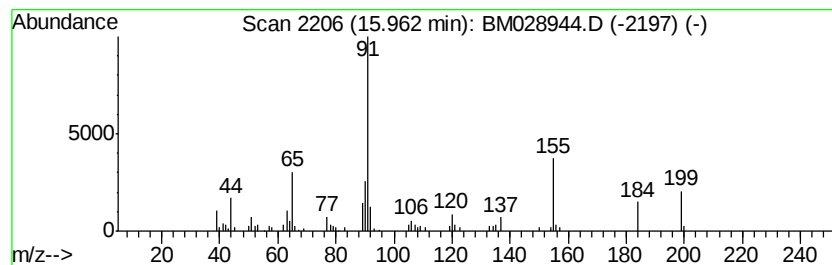
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Peak Number 3 Benzenesulfonamide, N-ethyl... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD		R.T.	
15.96	7.51 ng/ul	75523	Phenanthrene-d10		17.18	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzenesulfonamide, N-ethyl-2-me...	199	C9H13NO2S	001077-56-1	98
2		Benzenesulfonamide, N-ethyl-2-me...	199	C9H13NO2S	001077-56-1	49
3		4-Methyl-N-(1H-tetrazol-5-yl)-be...	239	C8H9N5O2S	006455-80-7	47
4		(O-p-Tosylisonitroso)malononitrile	249	C10H7N3O3S	020893-01-0	47
5		Benzenesulfonamide, N,4-dimethyl...	214	C8H10N2O3S	000080-11-5	47



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM022721\
Data File : BM028944.D
Acq On : 27 Feb 2021 06:53
Operator : CG/JU
Sample : M1498-06
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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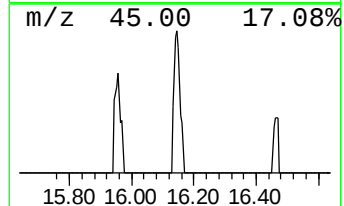
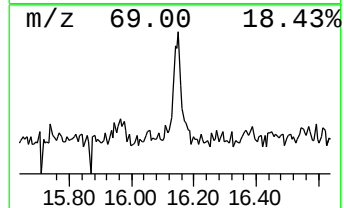
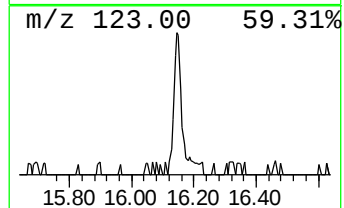
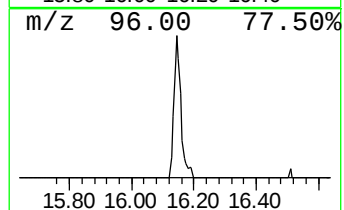
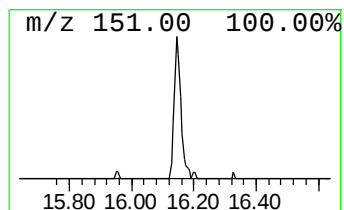
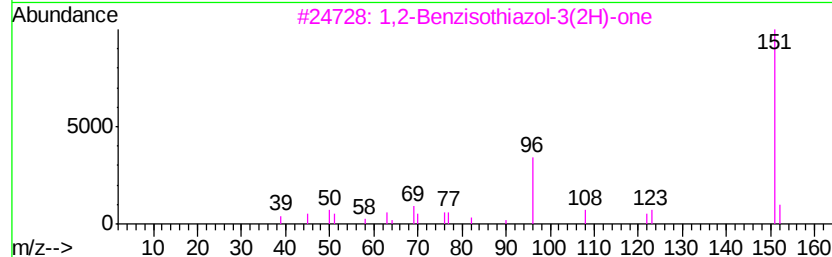
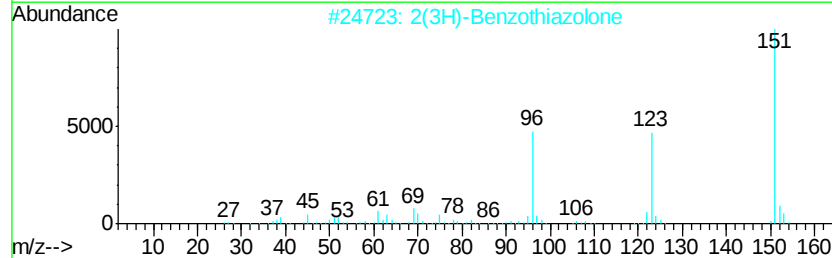
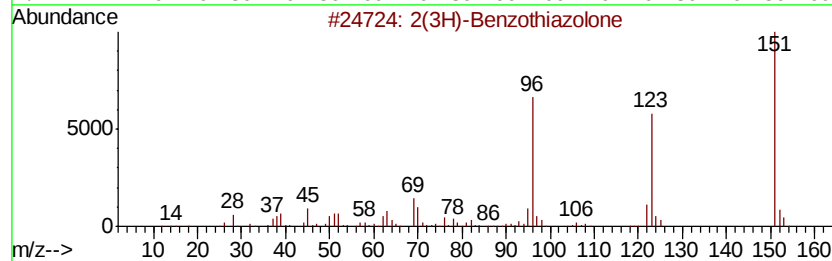
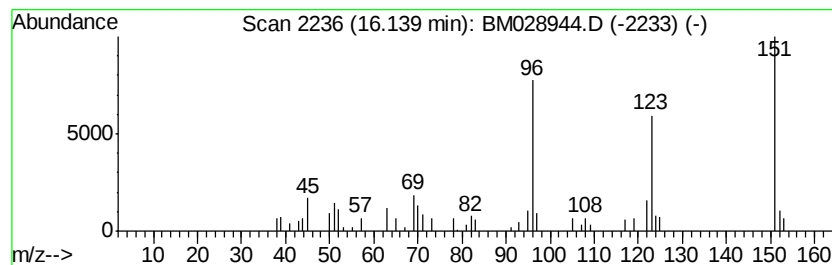
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 2(3H)-Benzothiazolone Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.14	2.96 ng/ul	29788	Phenanthrene-d10	17.18

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2(3H)-Benzothiazolone	151	C7H5NOS	000934-34-9	94
2		2(3H)-Benzothiazolone	151	C7H5NOS	000934-34-9	91
3		1,2-Benzisothiazol-3(2H)-one	151	C7H5NOS	002634-33-5	59
4		Thieno[2,3-b]pyridine-N-oxide	151	C7H5NOS	025557-50-0	46
5		Thieno[3,2-c]pyridin-4(5H)-one	151	C7H5NOS	027685-92-3	38



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM022721\
Data File : BM028944.D
Acq On : 27 Feb 2021 06:53
Operator : CG/JU
Sample : M1498-06
Misc :
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Instrument :
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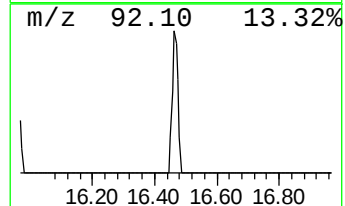
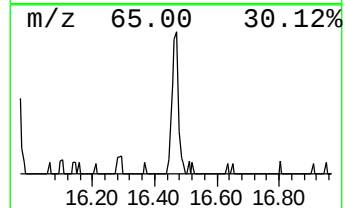
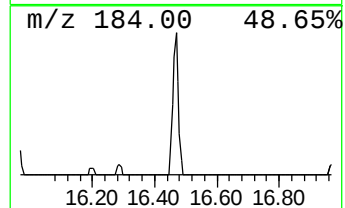
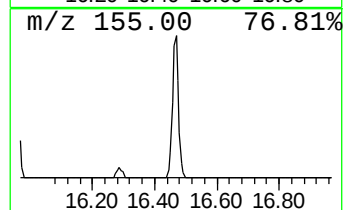
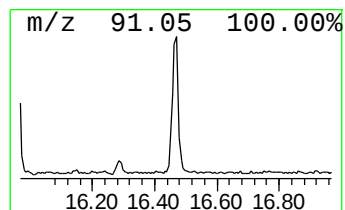
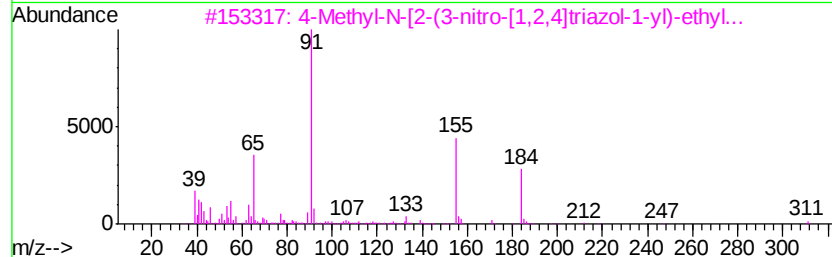
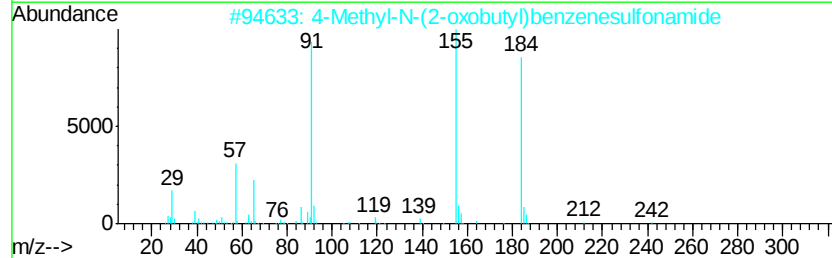
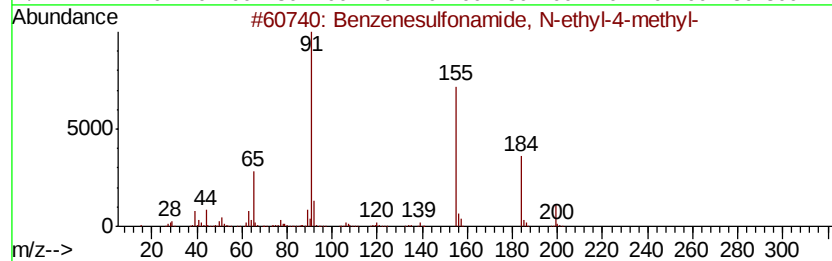
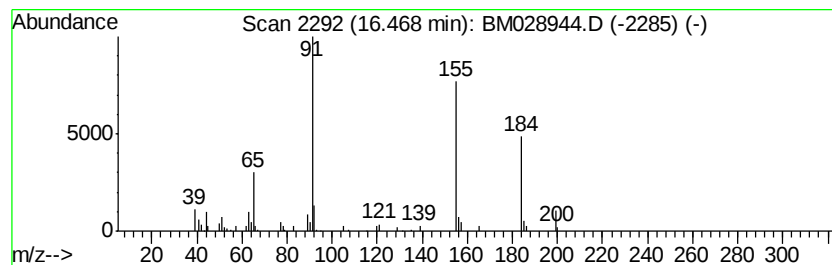
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Benzenesulfonamide, N-ethyl... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.47	4.08 ng/ul	41023	Phenanthrene-d10	17.18

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzenesulfonamide, N-ethyl-4-me...	199	C9H13NO2S	000080-39-7	96
2		4-Methyl-N-(2-oxobutyl)benzenesu...	241	C11H15NO3S	1000192-14-5	90
3		4-Methyl-N-[2-(3-nitro-[1,2,4]tr...	311	C11H13N5O4S	1000301-03-5	83
4		(Toluene-4-sulfonylamino)-acetic...	283	C13H17NO4S	1000190-48-0	78
5		Ethyl(E/Z)-.alpha.-cyano-4-nitro...	416	C19H16N2O7S	1000287-26-8	53



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM022721\
Data File : BM028944.D
Acq On : 27 Feb 2021 06:53
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Sample : M1498-06
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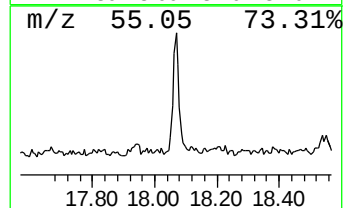
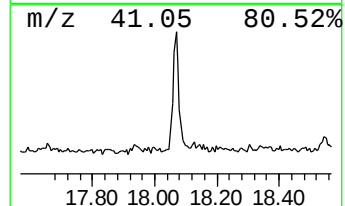
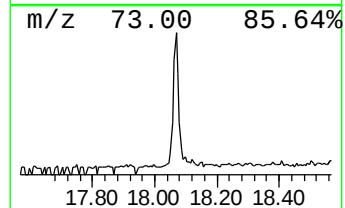
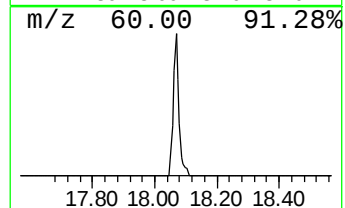
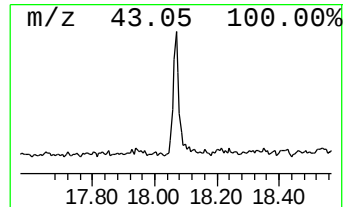
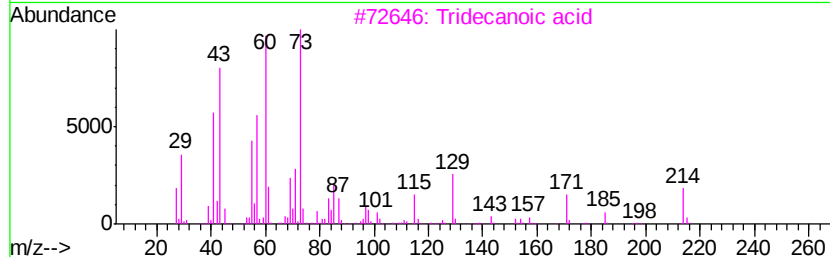
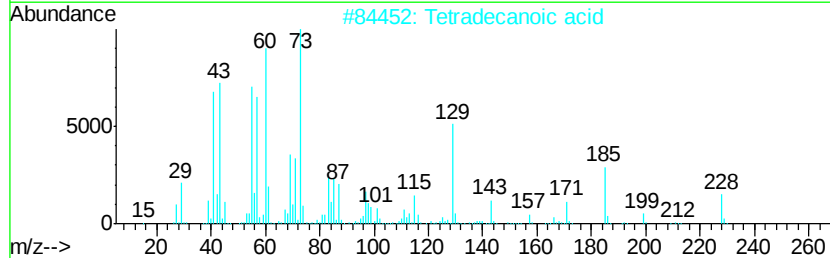
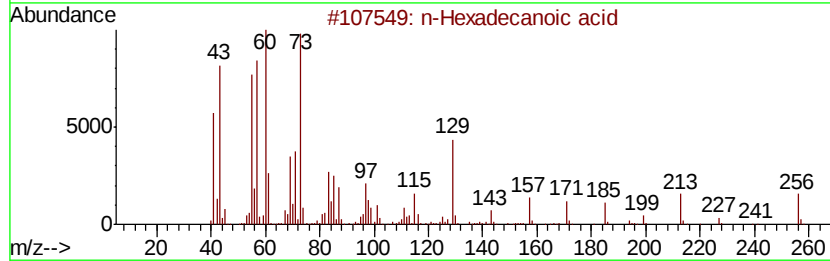
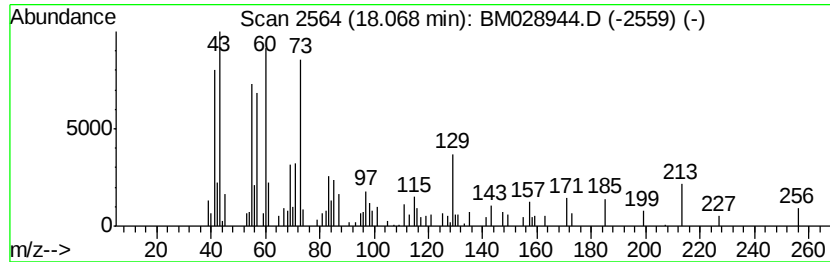
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.07	5.33 ng/ul	53654	Phenanthrene-d10	17.18

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2		Tetradecanoic acid	228	C14H28O2	000544-63-8	74
3		Tridecanoic acid	214	C13H26O2	000638-53-9	72
4		Pentadecanoic acid	242	C15H30O2	001002-84-2	72
5		Pentadecanoic acid	242	C15H30O2	001002-84-2	68



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM022721\
Data File : BM028944.D
Acq On : 27 Feb 2021 06:53
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Sample : M1498-06
Misc :
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Instrument :
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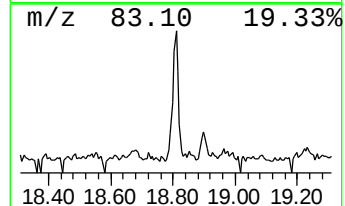
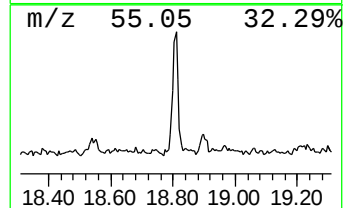
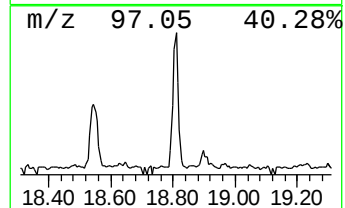
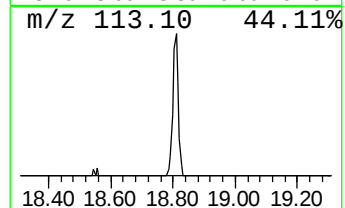
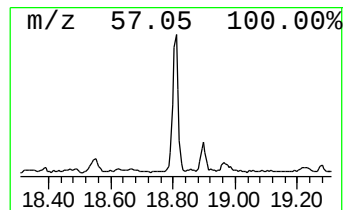
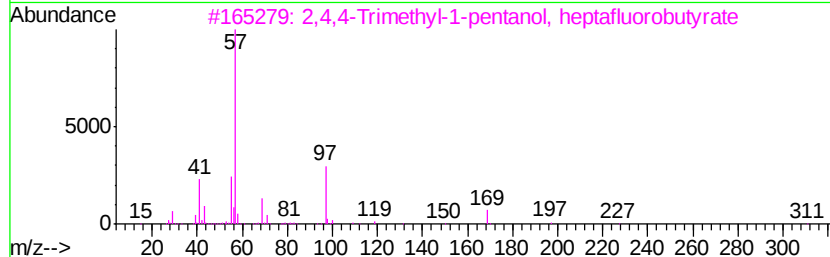
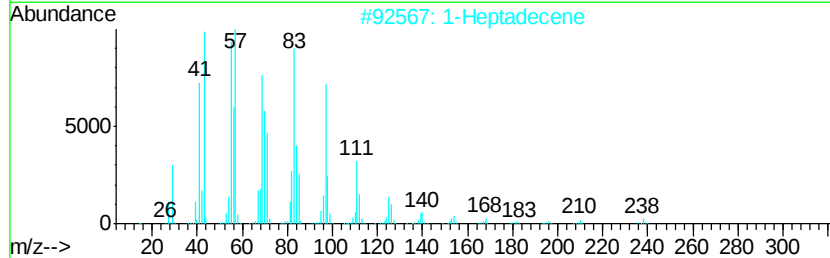
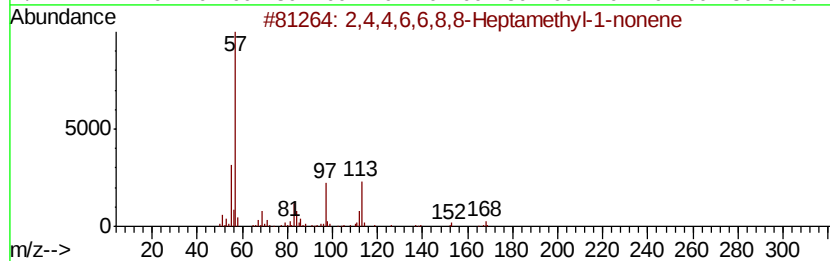
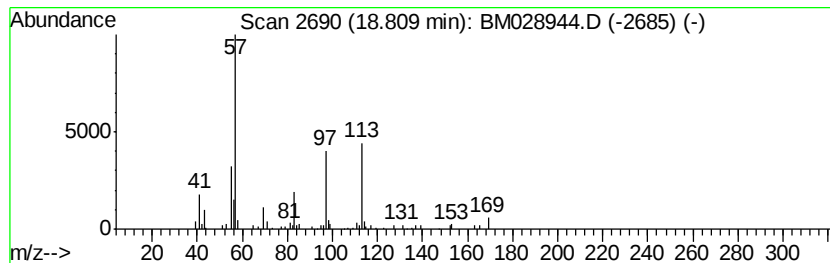
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.81	4.15 ng/ul	41790	Phenanthrene-d10	17.18

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2,4,4,6,6,8,8-Heptamethyl-1-nonene	224	C16H32	015796-04-0	47
2			1-Heptadecene	238	C17H34	006765-39-5	27
3			2,4,4-Trimethyl-1-pentanol, hept...	326	C12H17F7O2	1000365-01-4	25
4			Octane, 2-bromo-	192	C8H17Br	000557-35-7	25
5			Formic acid, 2,4,4-trimethylpent...	158	C9H18O2	1000368-95-2	25



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM022721\
Data File : BM028944.D
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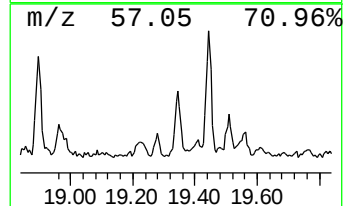
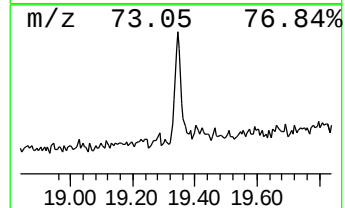
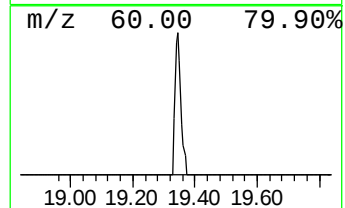
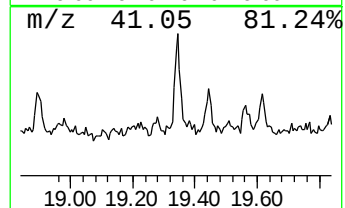
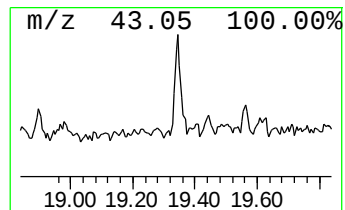
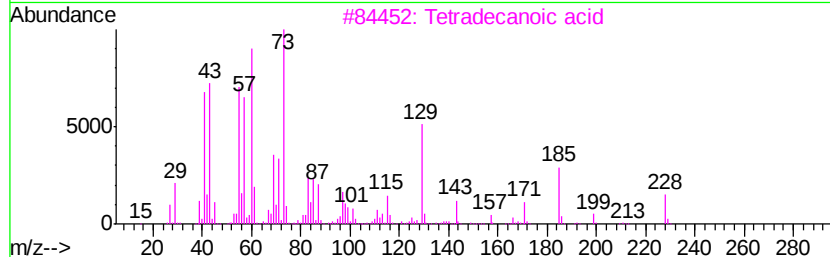
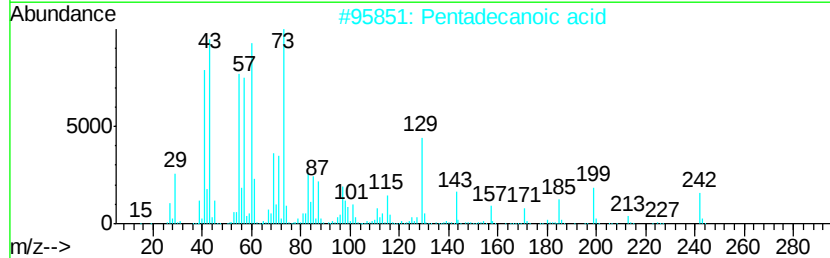
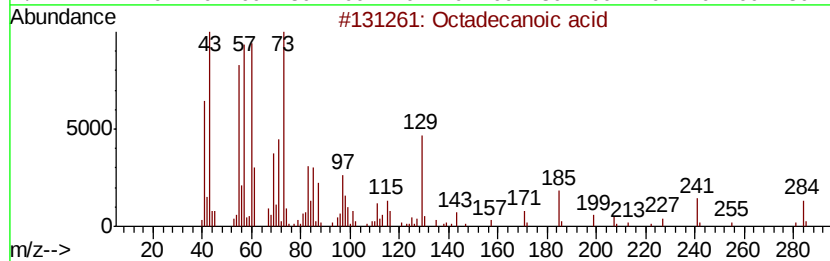
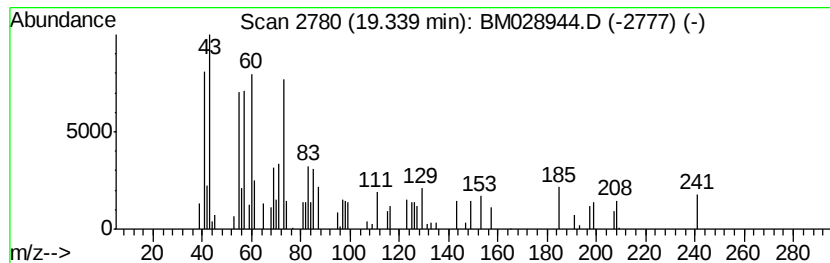
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Octadecanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.34	2.94 ng/ul	25533	Chrysene-d12	21.34

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecanoic acid	284	C18H36O2	000057-11-4	96
2		Pentadecanoic acid	242	C15H30O2	001002-84-2	62
3		Tetradecanoic acid	228	C14H28O2	000544-63-8	58
4		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	58
5		Tetradecanoic acid	228	C14H28O2	000544-63-8	53



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM022721\
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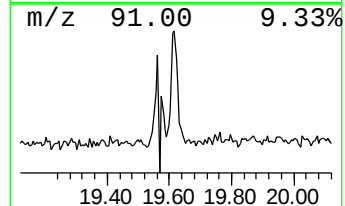
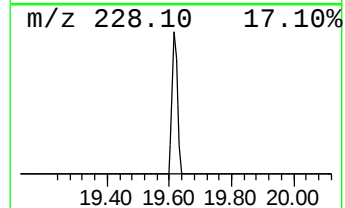
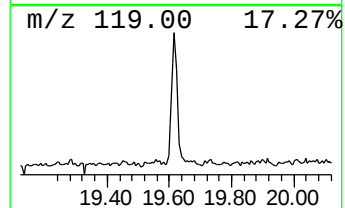
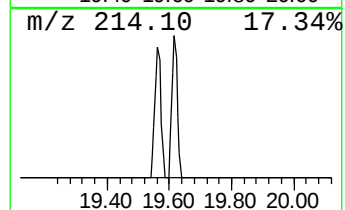
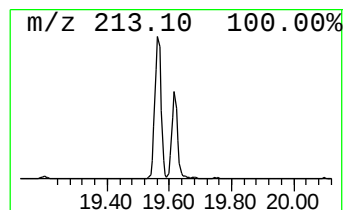
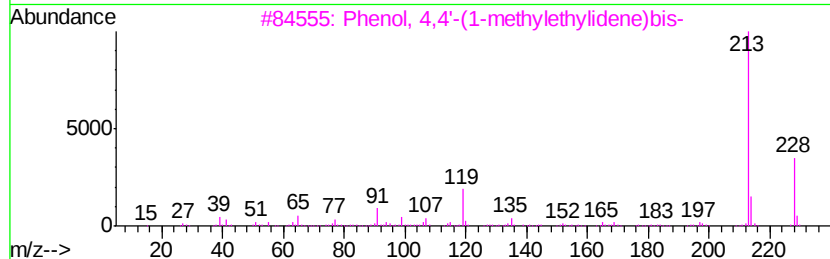
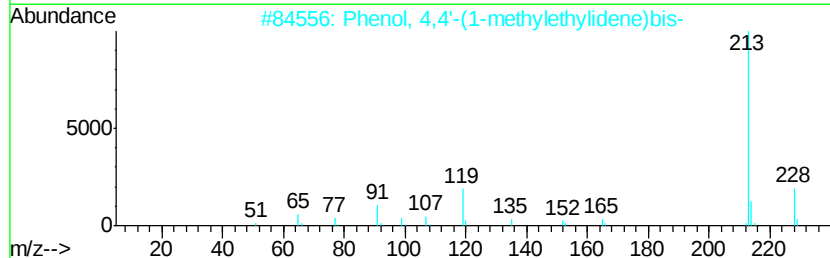
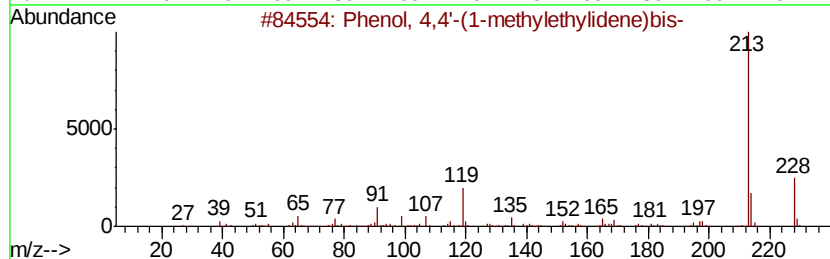
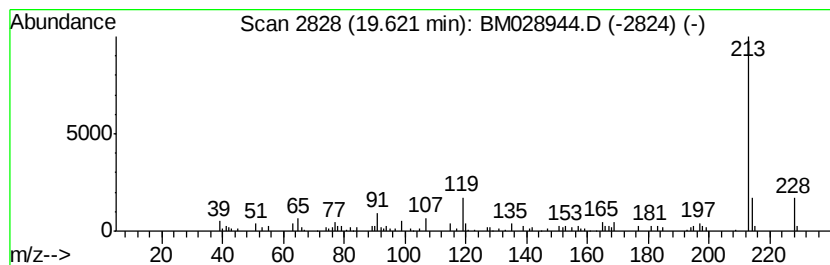
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Phenol, 4,4'-(1-methylethyl... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.62	6.16 ng/ul	53513	Chrysene-d12	21.34

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 4,4'-(1-methylethylidene)bis-	228	C15H16O2	000080-05-7	97	
2	Phenol, 4,4'-(1-methylethylidene)bis-	228	C15H16O2	000080-05-7	94	
3	Phenol, 4,4'-(1-methylethylidene)bis-	228	C15H16O2	000080-05-7	91	
4	2-Trimethylsilyl-3-trimethylsilyl-	228	C8H20N4Si2	1000373-05-5	64	
5	Carbanilic acid, p-phenyl-	213	C13H11NO2	004474-53-7	64	



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM022721\
Data File : BM028944.D
Acq On : 27 Feb 2021 06:53
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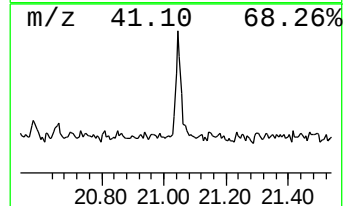
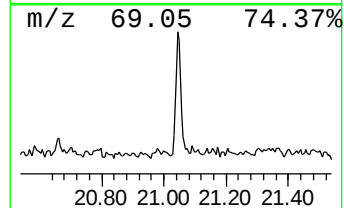
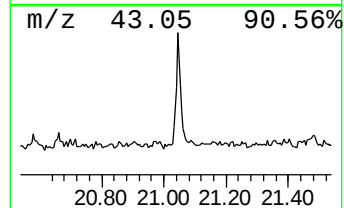
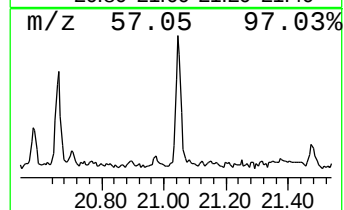
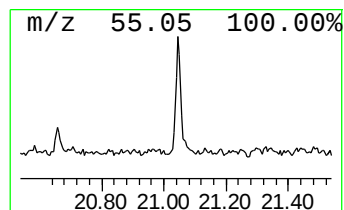
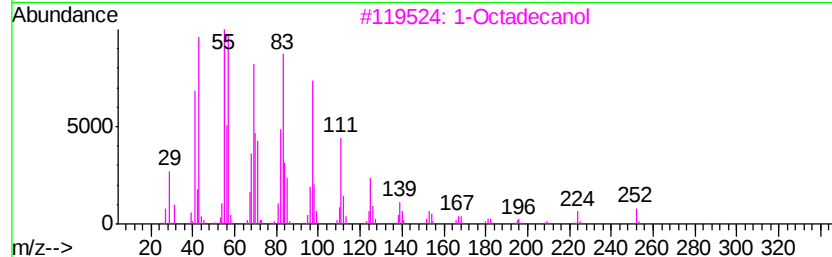
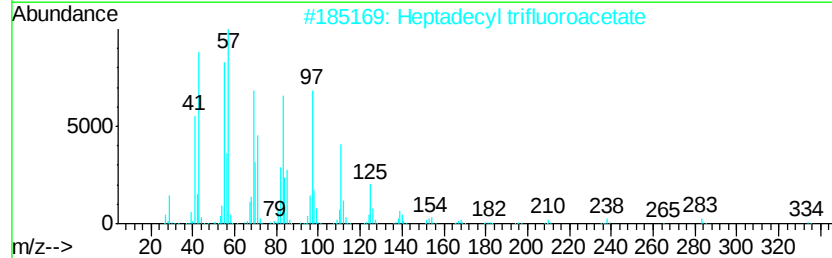
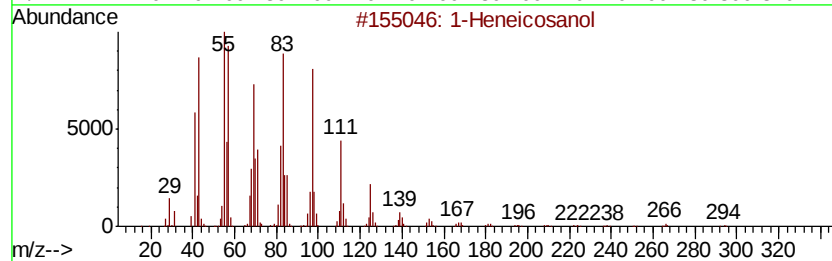
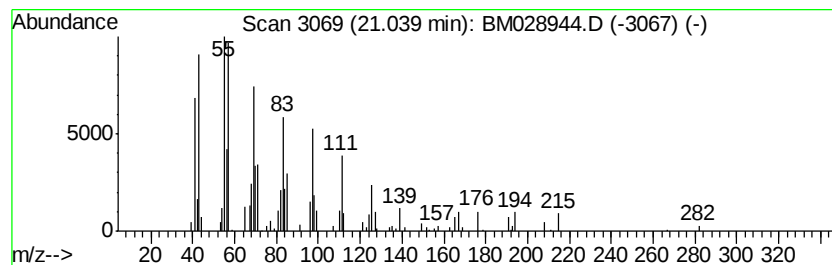
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM022721MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 10 1-Heneicosanol Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.04	4.55 ng/ul	39535	Chrysene-d12	21.34

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Heneicosanol	312	C21H44O	015594-90-8	94
2			Heptadecyl trifluoroacetate	352	C19H35F3O2	1000351-87-0	93
3			1-Octadecanol	270	C18H38O	000112-92-5	93
4			1-Nonadecene	266	C19H38	018435-45-5	90
5			Cycloeicosane	280	C20H40	000296-56-0	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM022721\
 Data File : BM028944.D
 Acq On : 27 Feb 2021 06:53
 Operator : CG/JU
 Sample : M1498-06
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 CB8F9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM022721MA.M
 Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzene, chloro-	5.05	9.1	ng/ul	45317	1	7.77	99856	20.0
unknown-01	13.32	2.3	ng/ul	21139	3	14.43	185628	20.0
Benzenesulfonamid...	15.96	7.5	ng/ul	75523	4	17.18	201260	20.0
2(3H)-Benzothiazo...	16.14	3.0	ng/ul	29788	4	17.18	201260	20.0
Benzenesulfonamid...	16.47	4.1	ng/ul	41023	4	17.18	201260	20.0
n-Hexadecanoic acid	18.07	5.3	ng/ul	53654	4	17.18	201260	20.0
(DEL) Alkane: Str...	18.81	4.2	ng/ul	41790	4	17.18	201260	20.0
Octadecanoic acid	19.34	2.9	ng/ul	25533	5	21.34	173699	20.0
Phenol, 4,4'-(1-m...	19.62	6.2	ng/ul	53513	5	21.34	173699	20.0
1-Heneicosanol	21.04	4.5	ng/ul	39535	5	21.34	173699	20.0