

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM022721\
 Data File : BM028946.D
 Acq On : 27 Feb 2021 08:05
 Operator : CG/JU
 Sample : M1498-02
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 CB8G1

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	25	rBV	5475	5735	1.39%	0.099%
2	3.157	25	29	38	rVB	21020	27946	6.76%	0.484%
3	4.346	227	231	238	rBV	21769	30905	7.47%	0.536%
4	5.040	345	349	357	rBB	26470	38068	9.21%	0.660%
5	6.869	656	660	665	rBB2	4619	5504	1.33%	0.095%
6	6.957	671	675	685	rVB2	22575	40200	9.72%	0.697%
7	7.110	696	701	709	rBB	60225	90391	21.86%	1.567%
8	7.304	728	734	742	rBV	76167	117360	28.38%	2.034%
9	7.769	808	813	822	rBB	73812	109364	26.45%	1.896%
10	7.863	823	829	835	rBB3	4216	6931	1.68%	0.120%
11	8.140	871	876	880	rBB	4846	6586	1.59%	0.114%
12	8.504	932	938	946	rBV	58479	91069	22.02%	1.579%
13	8.769	979	983	990	rBB	11244	16638	4.02%	0.288%
14	8.881	997	1002	1007	rBV2	4587	6860	1.66%	0.119%
15	8.945	1007	1013	1023	rVV	82982	129679	31.36%	2.248%
16	9.287	1065	1071	1079	rBB4	7009	11617	2.81%	0.201%
17	9.404	1086	1091	1097	rBB4	6265	10061	2.43%	0.174%
18	9.663	1129	1135	1145	rBB	69945	114177	27.61%	1.979%
19	10.110	1204	1211	1221	rVB3	5557	15317	3.70%	0.266%
20	10.204	1221	1227	1237	rBV	102794	168196	40.67%	2.915%
21	10.522	1276	1281	1284	rVV	6064	9719	2.35%	0.168%
22	10.569	1284	1289	1297	rVV	96777	155769	37.67%	2.700%
23	10.722	1306	1315	1326	rBV	83428	146730	35.48%	2.543%
24	11.392	1420	1429	1432	rBV3	2492	5152	1.25%	0.089%
25	11.422	1432	1434	1440	rVV3	3563	5402	1.31%	0.094%
26	11.786	1491	1496	1499	rVV3	2512	4158	1.01%	0.072%
27	11.928	1512	1520	1525	rBV2	11032	18434	4.46%	0.320%
28	12.098	1543	1549	1554	rBV2	4487	8656	2.09%	0.150%
29	12.310	1578	1585	1588	rBV3	2163	3980	0.96%	0.069%
30	12.586	1628	1632	1637	rBV4	4510	6678	1.61%	0.116%
31	13.022	1702	1706	1715	rVB4	2603	5917	1.43%	0.103%
32	13.239	1737	1743	1747	rBV5	3978	7108	1.72%	0.123%
33	13.322	1754	1757	1763	rVB2	6323	8328	2.01%	0.144%
34	13.392	1765	1769	1774	rVB3	3065	4301	1.04%	0.075%

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35	13.486	1781	1785	1789	rBV3	4224	5775	1.40%	0.100%
36	13.622	1805	1808	1812	rBV5	2230	3942	0.95%	0.068%
37	13.669	1812	1816	1822	rVB6	3952	6048	1.46%	0.105%
38	13.851	1839	1847	1852	rVV	173481	237353	57.39%	4.114%
39	13.898	1852	1855	1859	rVB	14189	16539	4.00%	0.287%
40	14.004	1867	1873	1874	rBV6	2562	4491	1.09%	0.078%
41	14.080	1882	1886	1889	rBV5	5570	7383	1.79%	0.128%
42	14.127	1889	1894	1900	rVV	198081	281940	68.18%	4.887%
43	14.263	1913	1917	1923	rBV	9054	14535	3.51%	0.252%
44	14.363	1929	1934	1939	rBV3	17085	24805	6.00%	0.430%
45	14.433	1940	1946	1953	rVB	135193	195793	47.34%	3.394%
46	14.498	1953	1957	1961	rBV2	8047	11013	2.66%	0.191%
47	14.692	1985	1990	1999	rBV	14599	32737	7.92%	0.567%
48	14.874	2017	2021	2024	rVB3	4319	5948	1.44%	0.103%
49	15.051	2046	2051	2059	rBV10	5424	12092	2.92%	0.210%
50	15.186	2069	2074	2083	rBV	51799	83110	20.10%	1.441%
51	15.263	2083	2087	2092	rVB	7035	11238	2.72%	0.195%
52	15.316	2092	2096	2099	rBV2	5460	6978	1.69%	0.121%
53	15.345	2099	2101	2106	rVB4	3522	4473	1.08%	0.078%
54	15.427	2109	2115	2121	rBV	252704	353852	85.57%	6.134%
55	15.486	2121	2125	2129	rVB4	6046	9880	2.39%	0.171%
56	15.574	2132	2140	2149	rBV	62178	94210	22.78%	1.633%
57	15.663	2151	2155	2158	rBV	10599	14254	3.45%	0.247%
58	15.692	2158	2160	2162	rVB3	5110	4002	0.97%	0.069%
59	15.851	2184	2187	2191	rVB3	5830	7112	1.72%	0.123%
60	15.892	2191	2194	2198	rBV5	2584	3927	0.95%	0.068%
61	15.957	2199	2205	2215	rBV	92518	129068	31.21%	2.237%
62	16.110	2226	2231	2235	rVV3	6472	10429	2.52%	0.181%
63	16.151	2235	2238	2243	rVB5	4219	5942	1.44%	0.103%
64	16.286	2257	2261	2264	rBV3	4567	7411	1.79%	0.128%
65	16.468	2287	2292	2296	rBV	46349	62008	14.99%	1.075%
66	16.510	2297	2299	2308	rVB4	5711	7656	1.85%	0.133%
67	16.598	2310	2314	2317	rBV6	2712	4000	0.97%	0.069%
68	16.733	2333	2337	2344	rBV	11212	17012	4.11%	0.295%
69	17.039	2385	2389	2391	rBV5	3019	4639	1.12%	0.080%
70	17.174	2407	2412	2419	rVV	147975	209449	50.65%	3.631%
71	17.274	2424	2429	2435	rBV	268664	362564	87.67%	6.285%

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72	17.404	2447	2451	2456	rBV5	3219	5594	1.35%	0.097%
73	17.468	2456	2462	2468	rBV8	3837	9535	2.31%	0.165%
74	17.939	2538	2542	2546	rBV4	3545	5920	1.43%	0.103%
75	18.068	2559	2564	2571	rBV	94452	132082	31.94%	2.289%
76	18.209	2585	2588	2591	rBV4	2699	4003	0.97%	0.069%
77	18.545	2641	2645	2651	rVB	6938	11372	2.75%	0.197%
78	18.645	2660	2662	2673	rVB8	5100	8467	2.05%	0.147%
79	18.809	2685	2690	2694	rBV	45863	61397	14.85%	1.064%
80	18.845	2694	2696	2701	rVB4	11998	15159	3.67%	0.263%
81	18.898	2701	2705	2709	rVB	12304	14027	3.39%	0.243%
82	18.962	2713	2716	2722	rVB5	3441	6688	1.62%	0.116%
83	19.198	2753	2756	2758	rBV3	6884	9251	2.24%	0.160%
84	19.280	2767	2770	2774	rVB2	6999	8087	1.96%	0.140%
85	19.345	2777	2781	2788	rBV	60802	79369	19.19%	1.376%
86	19.445	2794	2798	2802	rVB2	12086	14855	3.59%	0.257%
87	19.562	2812	2818	2823	rBV	299048	396549	95.89%	6.874%
88	19.615	2823	2827	2835	rVB	160183	210806	50.98%	3.654%
89	19.974	2884	2888	2893	rBV3	5833	8759	2.12%	0.152%
90	20.450	2967	2969	2973	rVB	5078	4407	1.07%	0.076%
91	20.580	2988	2991	2995	rBV3	11503	13978	3.38%	0.242%
92	20.662	3001	3005	3009	rBV	10539	13613	3.29%	0.236%
93	21.045	3066	3070	3077	rBV	65624	76285	18.45%	1.322%
94	21.339	3115	3120	3127	rBV	177910	223334	54.00%	3.871%
95	21.897	3213	3215	3218	rBV2	10310	11293	2.73%	0.196%
96	22.350	3289	3292	3300	rVB3	16572	25746	6.23%	0.446%
97	22.844	3373	3376	3381	rVB	21990	28269	6.84%	0.490%
98	23.397	3464	3470	3477	rBV	239204	413547	100.00%	7.168%
99	23.539	3488	3494	3502	rVB2	113552	208717	50.47%	3.618%
100	24.021	3572	3576	3581	rVB	24510	41403	10.01%	0.718%

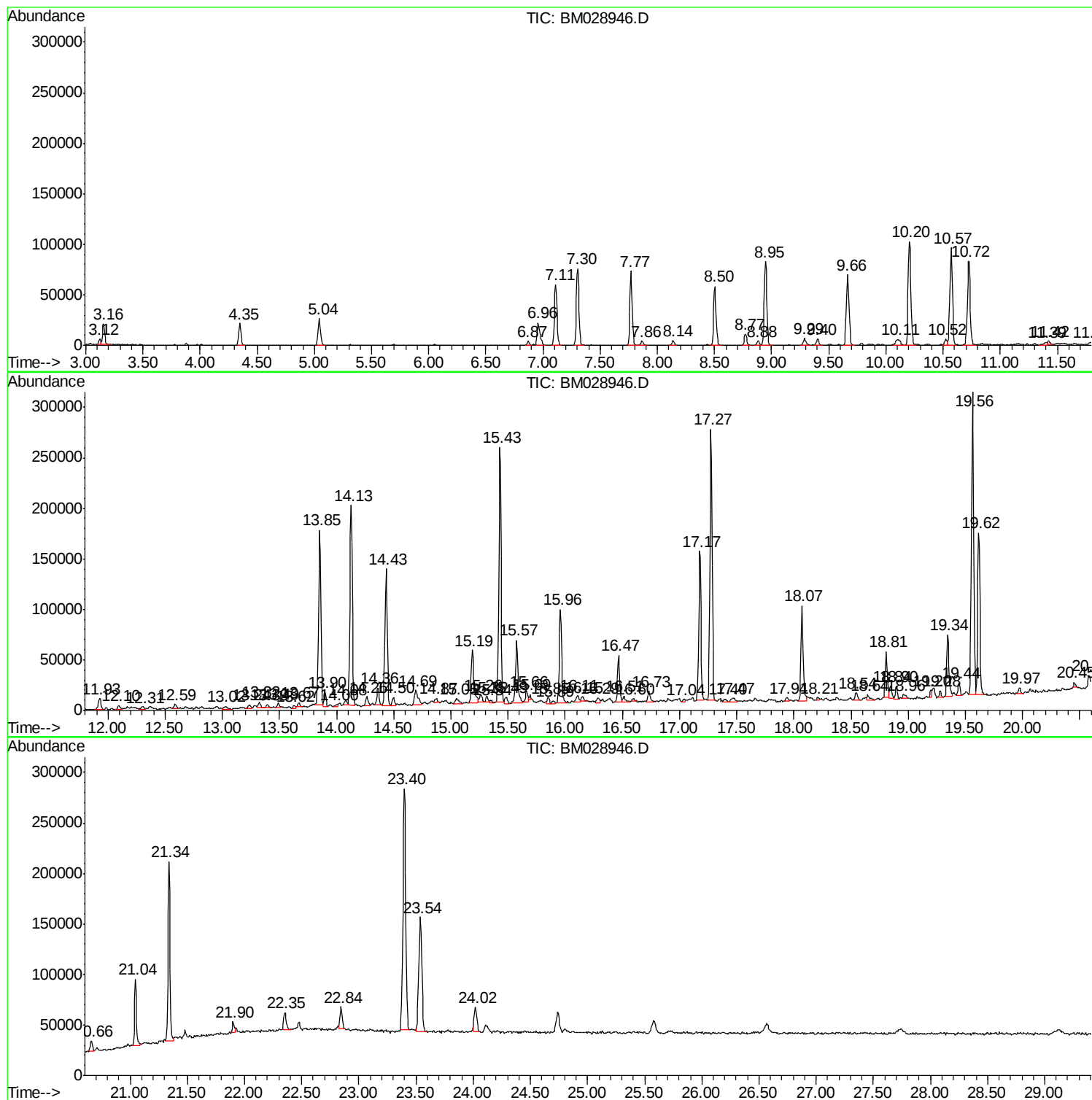
Sum of corrected areas: 5769056

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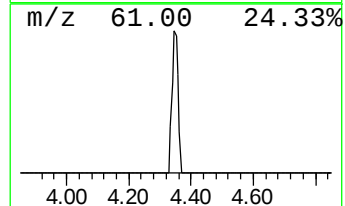
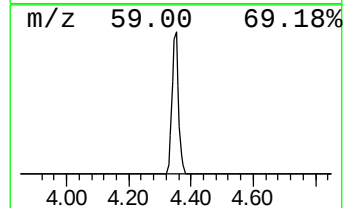
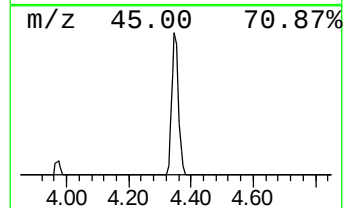
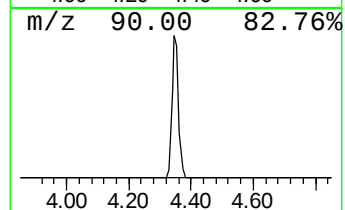
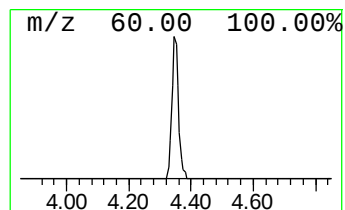
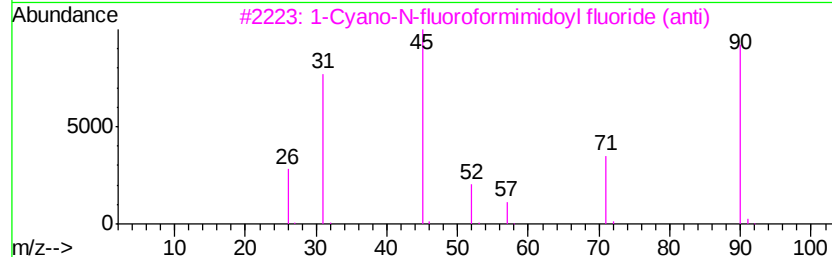
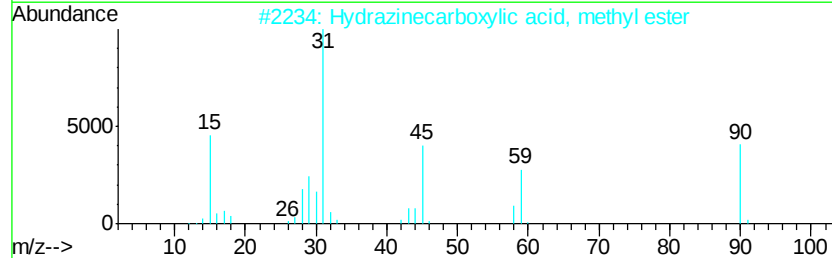
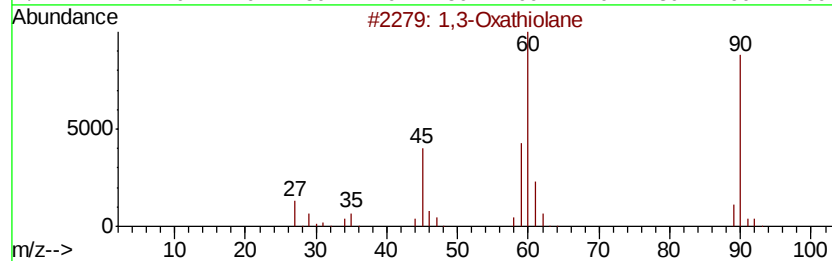
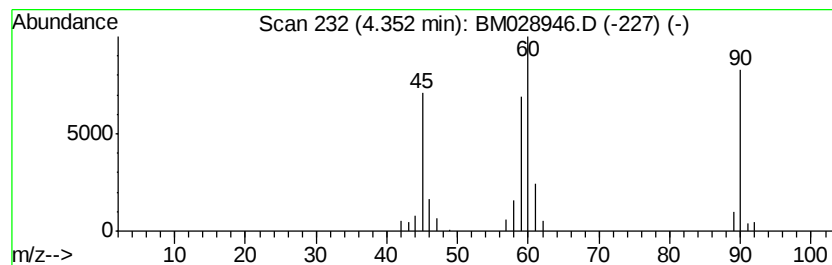
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM022721MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 1,3-Oxathiolane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.35	5.65 ng/ul	30905	1,4-Dichlorobenzene-d4	7.77

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,3-Oxathiolane	90	C3H6OS	002094-97-5	87
2			Hydrazinecarboxylic acid, methyl...	90	C2H6N2O2	006294-89-9	59
3			1-Cyano-N-fluoroformimidovl fluo...	90	C2F2N2	014011-05-3	40
4			3-Thietanol	90	C3H6OS	010304-16-2	40
5			Hydrazinecarboxylic acid, methyl...	90	C2H6N2O2	006294-89-9	38



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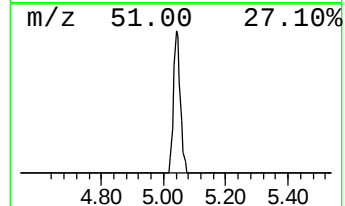
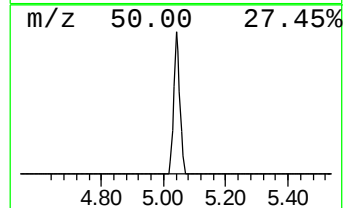
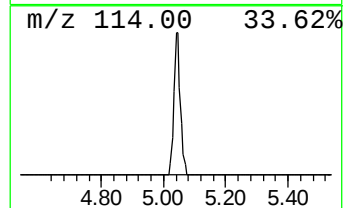
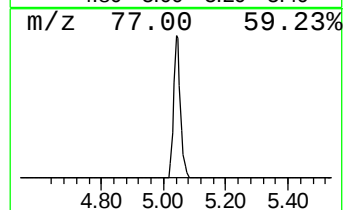
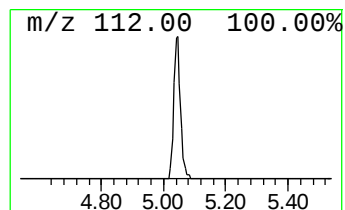
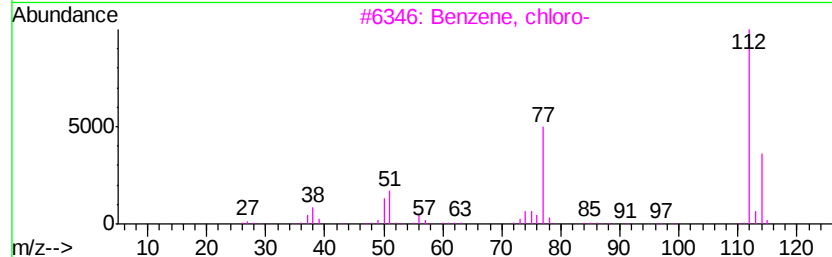
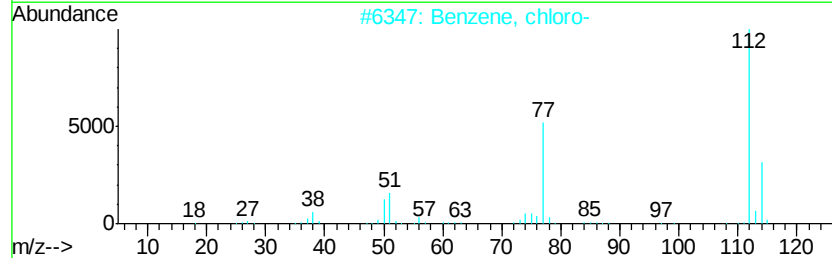
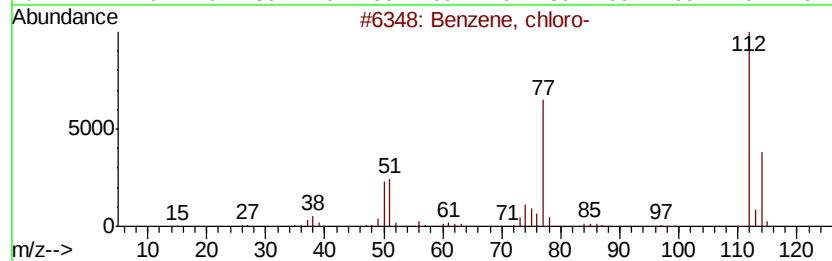
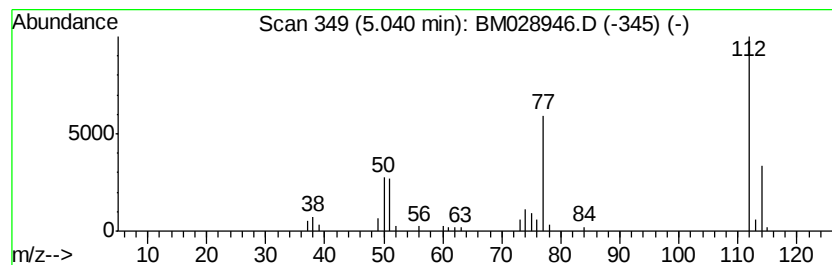
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Benzene, chloro- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.04	6.96 ng/ul	38068	1,4-Dichlorobenzene-d4	7.77

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, chloro-	112	C6H5Cl	000108-90-7	94
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	86
5			Phenol, 3-fluoro-	112	C6H5FO	000372-20-3	16



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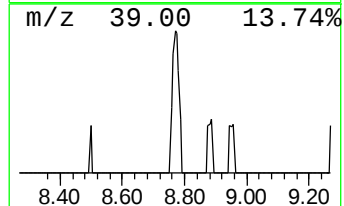
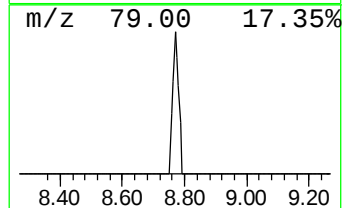
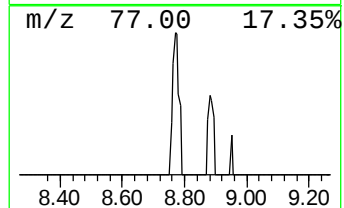
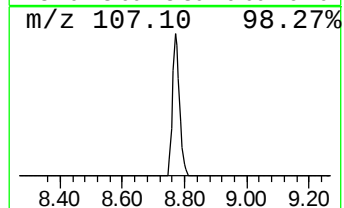
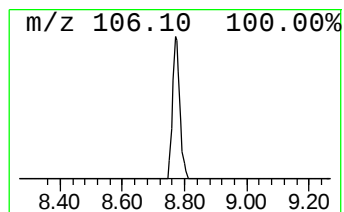
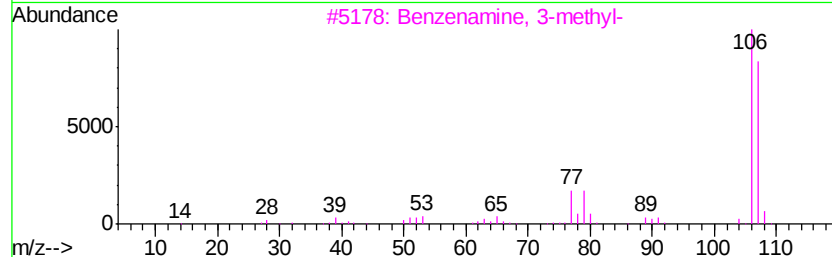
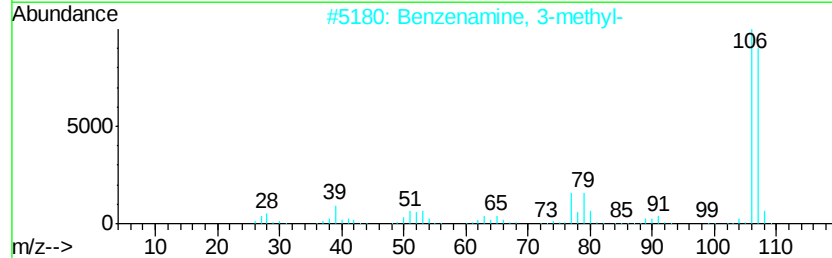
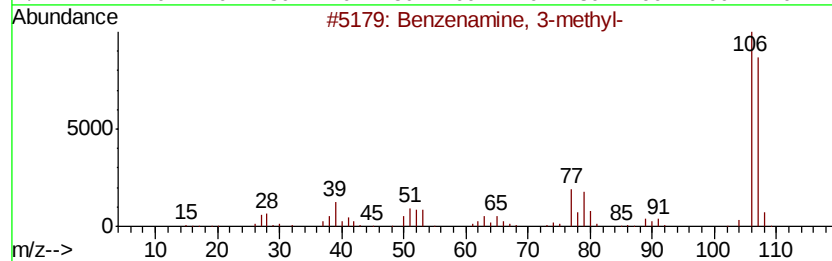
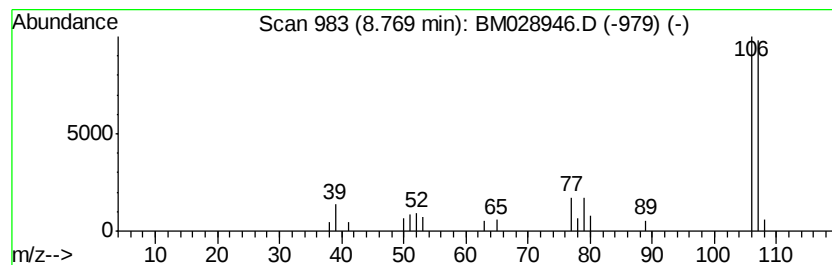
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Peak Number 3 Benzenamine, 3-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.77	3.04 ng/ul	16638	1,4-Dichlorobenzene-d4	7.77

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzenamine, 3-methyl-	107	C7H9N	000108-44-1	91
2		Benzenamine, 3-methyl-	107	C7H9N	000108-44-1	91
3		Benzenamine, 3-methyl-	107	C7H9N	000108-44-1	91
4		Benzenamine, 3-methyl-	107	C7H9N	000108-44-1	91
5		Benzenamine, 3-methyl-	107	C7H9N	000108-44-1	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM022721\
Data File : BM028946.D
Acq On : 27 Feb 2021 08:05
Operator : CG/JU
Sample : M1498-02
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
CB8G1

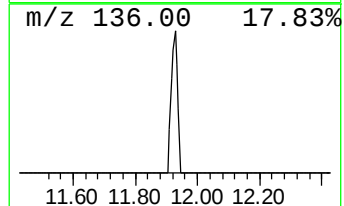
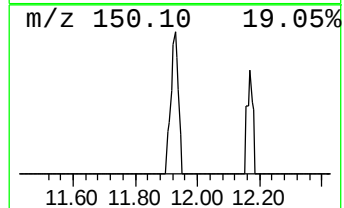
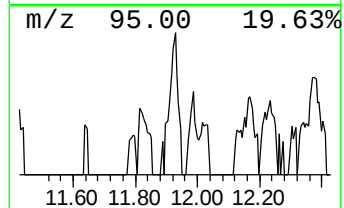
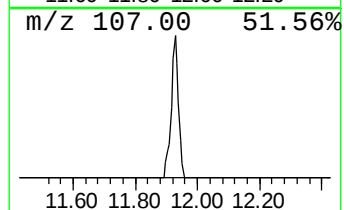
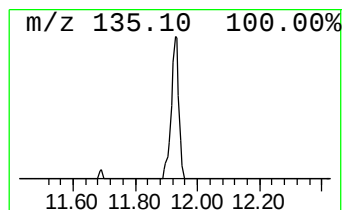
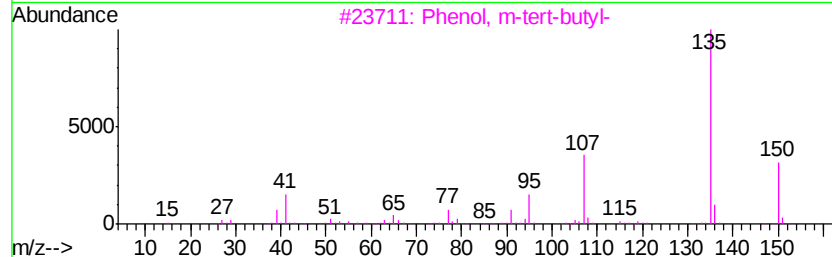
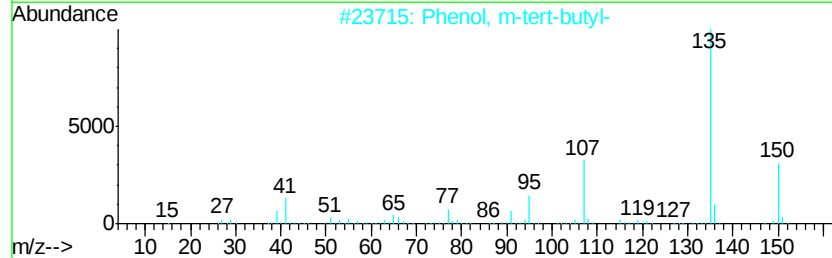
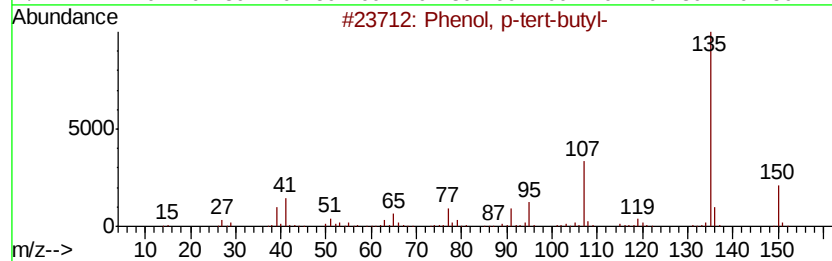
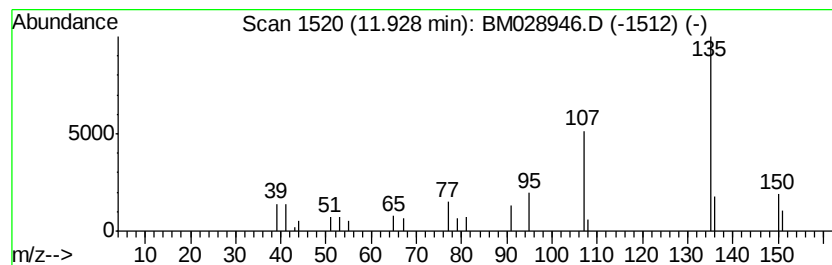
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Phenol, p-tert-butyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.93	2.37 ng/ul	18434	Naphthalene-d8	10.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	90
2		Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	83
3		Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	78
4		Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	72
5		Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	72



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM022721\
Data File : BM028946.D
Acq On : 27 Feb 2021 08:05
Operator : CG/JU
Sample : M1498-02
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
CB8G1

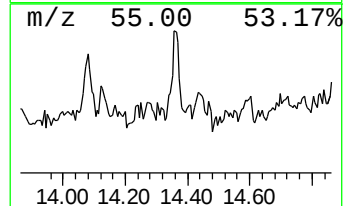
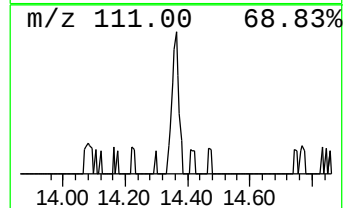
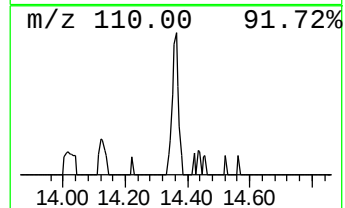
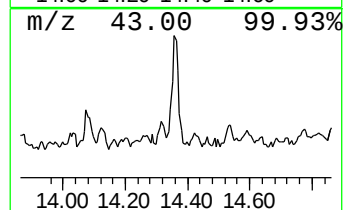
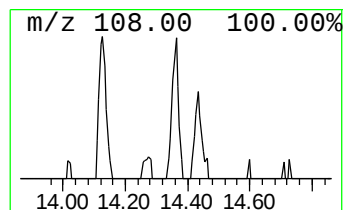
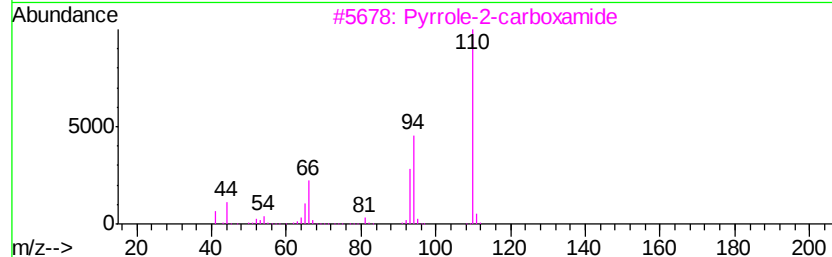
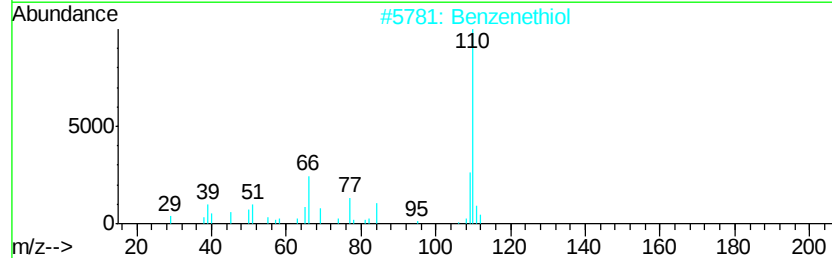
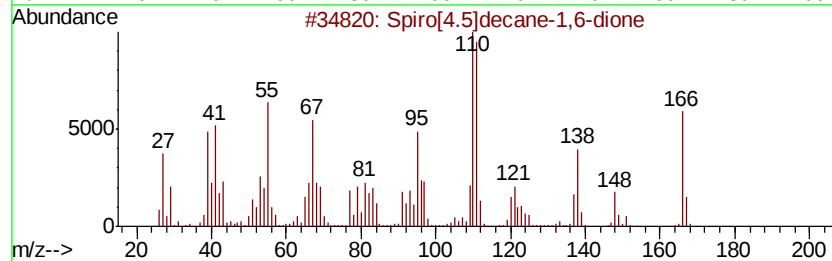
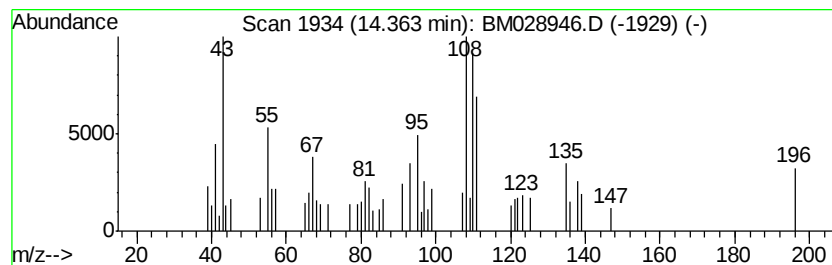
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 unknown-01 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.36	2.53 ng/ul	24805	Acenaphthene-d10	14.43

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Spino[4.5]decane-1,6-dione	166	C10H14O2	036803-48-2	43
2		Benzenethiol	110	C6H6S	000108-98-5	30
3		Pyrrole-2-carboxamide	110	C5H6N2O	004551-72-8	27
4		Benzenethiol	110	C6H6S	000108-98-5	27
5		Bicyclo[2.2.1]heptane, exo-2-met...	168	C10H16O2	1000138-90-3	22



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM022721\
Data File : BM028946.D
Acq On : 27 Feb 2021 08:05
Operator : CG/JU
Sample : M1498-02
Misc :
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Instrument :
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ClientSampleId :
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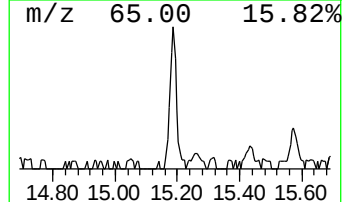
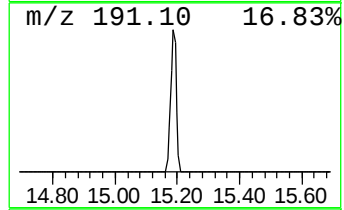
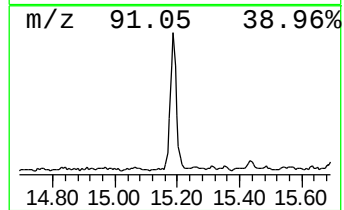
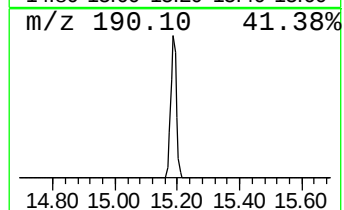
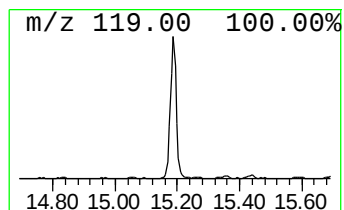
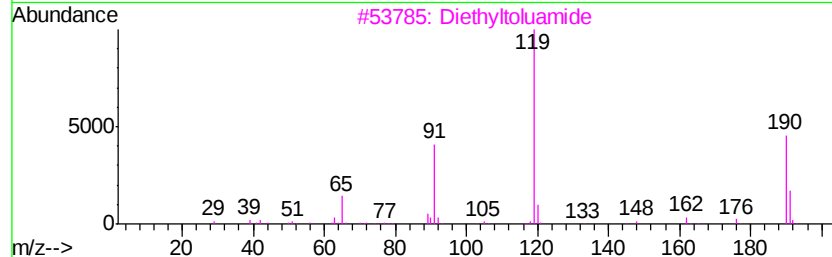
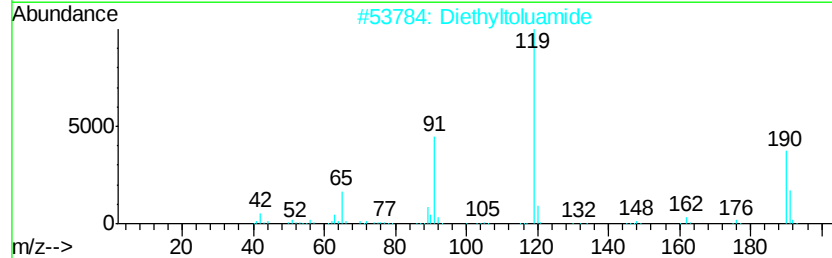
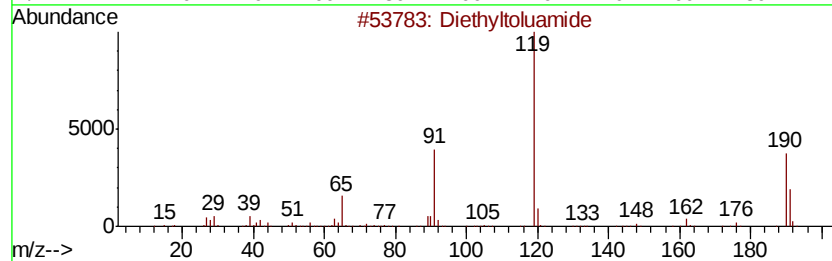
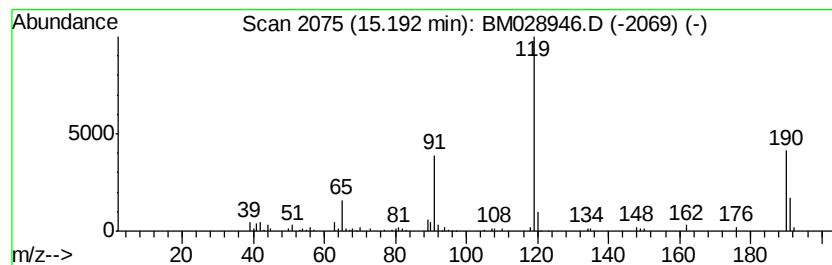
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Diethyltoluamide Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.19	8.49 ng/ul	83110	Acenaphthene-d10	14.43

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Diethyltoluamide	191	C12H17NO	000134-62-3	96
2		Diethyltoluamide	191	C12H17NO	000134-62-3	96
3		Diethyltoluamide	191	C12H17NO	000134-62-3	96
4		Benzamide, N,N-diethyl-4-methyl-	191	C12H17NO	002728-05-4	91
5		Diethyltoluamide	191	C12H17NO	000134-62-3	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM022721\
Data File : BM028946.D
Acq On : 27 Feb 2021 08:05
Operator : CG/JU
Sample : M1498-02
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
CB8G1

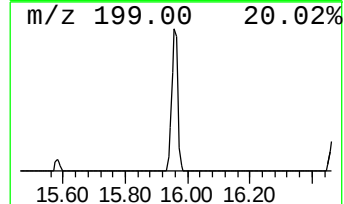
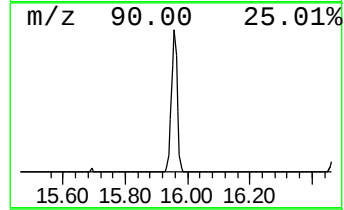
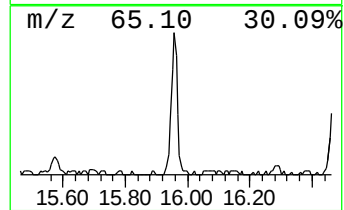
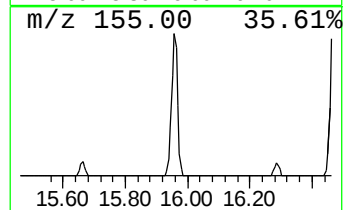
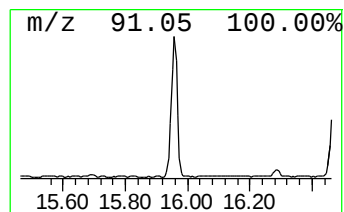
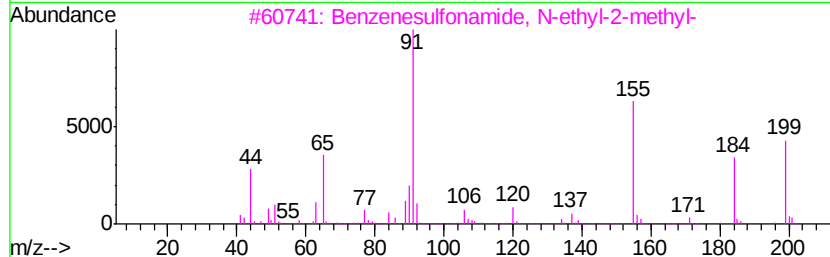
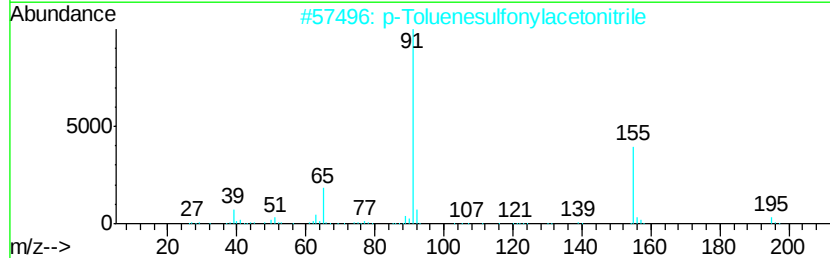
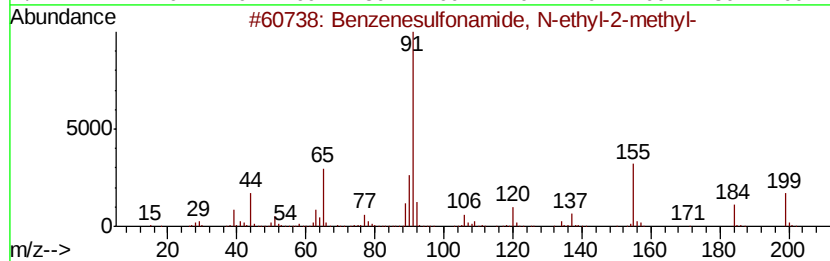
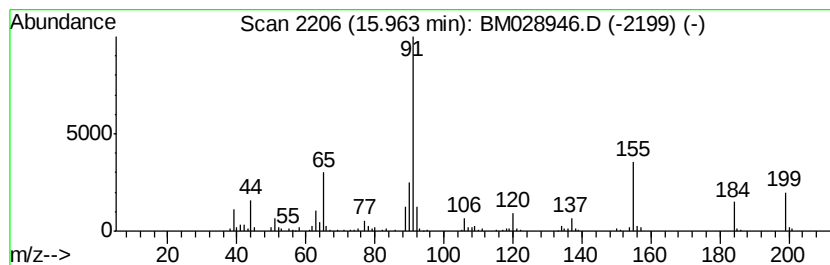
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Benzenesulfonamide, N-ethyl... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.96	12.32 ng/ul	129068	Phenanthrene-d10	17.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzenesulfonamide, N-ethyl-2-me...	199	C9H13NO2S	001077-56-1	99
2		p-Toluenesulfonylacetonitrile	195	C9H9NO2S	005697-44-9	47
3		Benzenesulfonamide, N-ethyl-2-me...	199	C9H13NO2S	001077-56-1	47
4		p-Toluenesulfonylacetonitrile	195	C9H9NO2S	005697-44-9	47
5		Benzenesulfonic acid, 4-methyl-,...	200	C9H12O3S	000080-40-0	46



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM022721\
Data File : BM028946.D
Acq On : 27 Feb 2021 08:05
Operator : CG/JU
Sample : M1498-02
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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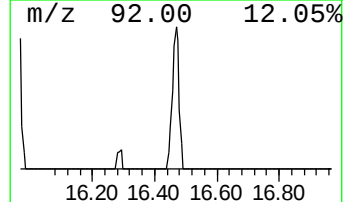
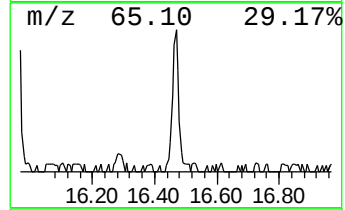
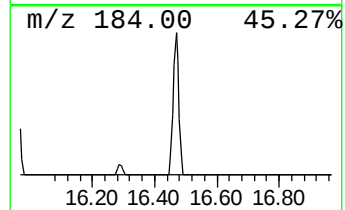
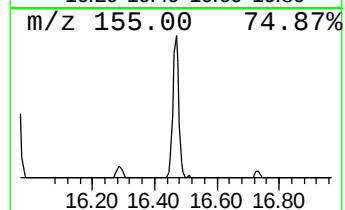
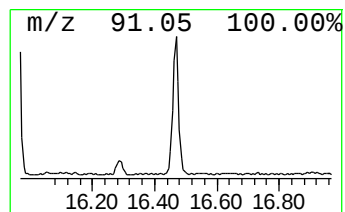
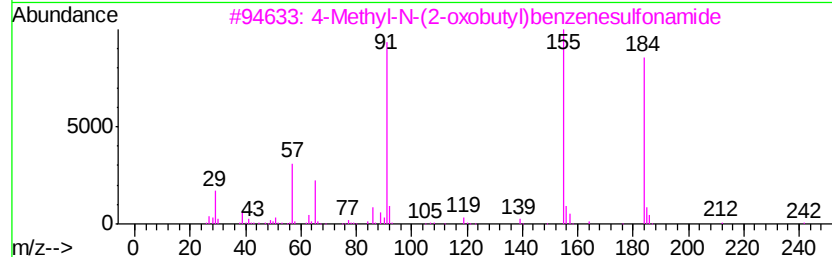
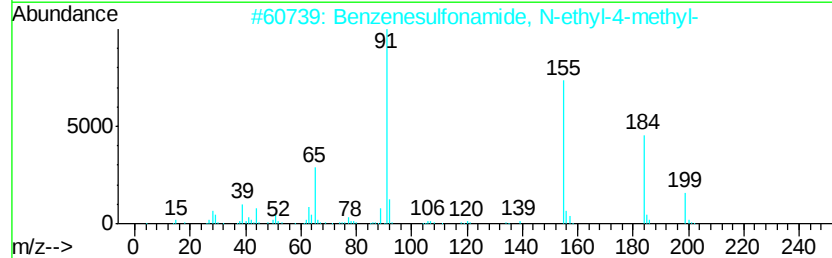
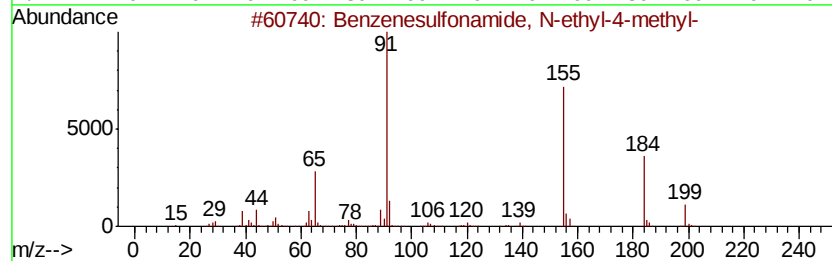
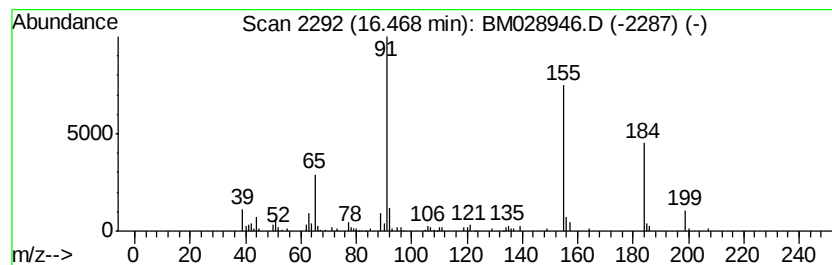
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Benzenesulfonamide, N-ethyl... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.47	5.92 ng/ul	62008	Phenanthrene-d10	17.17

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzenesulfonamide, N-ethyl-4-me...	199	C9H13NO2S	000080-39-7	98
2			Benzenesulfonamide, N-ethyl-4-me...	199	C9H13NO2S	000080-39-7	95
3			4-Methyl-N-(2-oxobutyl)benzenesu...	241	C11H15NO3S	1000192-14-5	86
4			4-Methyl-N-[2-(3-nitro-[1,2,4]tr...	311	C11H13N5O4S	1000301-03-5	83
5			4-Toluenesulfonylmethyl isocyanide	195	C9H9NO2S	036635-61-7	58



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM022721\
Data File : BM028946.D
Acq On : 27 Feb 2021 08:05
Operator : CG/JU
Sample : M1498-02
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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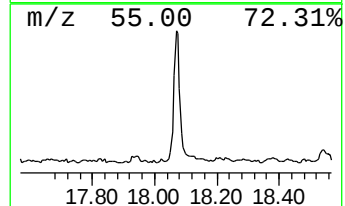
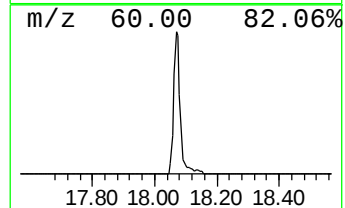
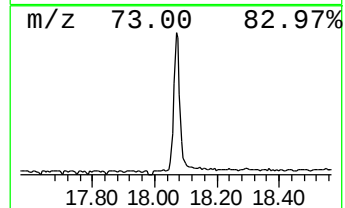
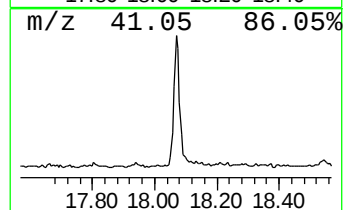
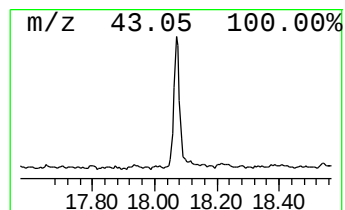
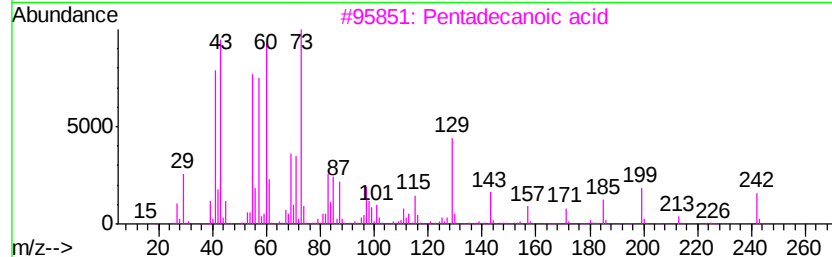
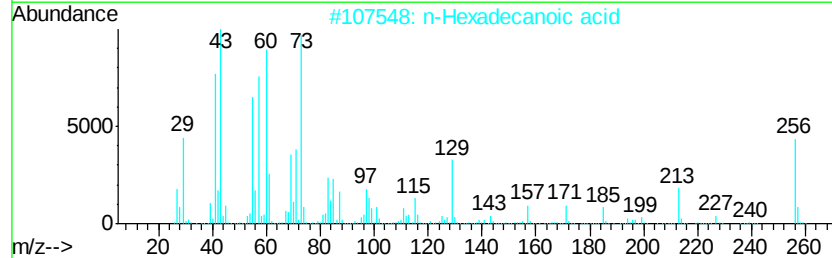
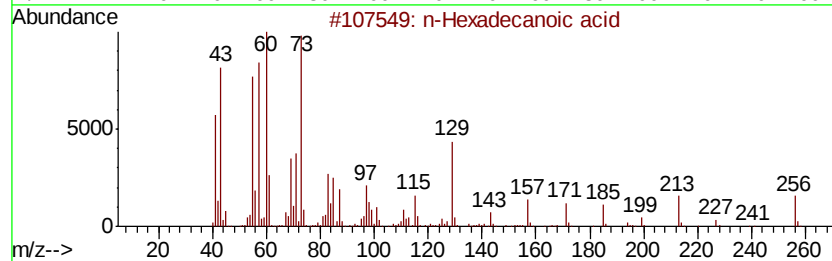
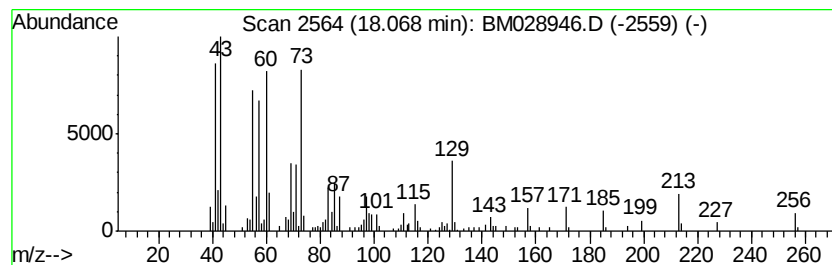
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.07	12.61 ng/ul	132082	Phenanthrene-d10	17.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	93
3		Pentadecanoic acid	242	C15H30O2	001002-84-2	90
4		Tridecanoic acid	214	C13H26O2	000638-53-9	76
5		Tridecanoic acid	214	C13H26O2	000638-53-9	76



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM022721\
Data File : BM028946.D
Acq On : 27 Feb 2021 08:05
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Sample : M1498-02
Misc :
ALS Vial : 35 Sample Multiplier: 1

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ClientSampleId :
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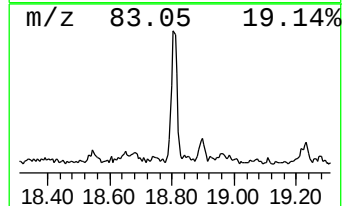
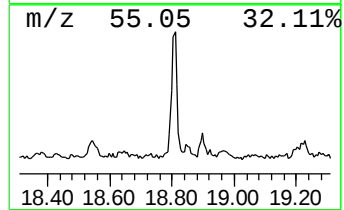
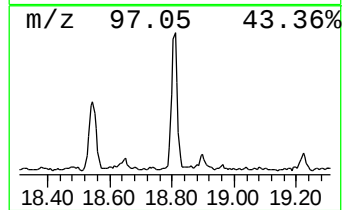
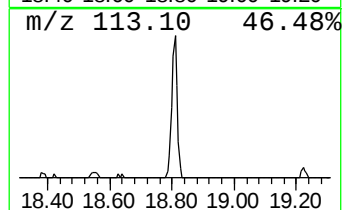
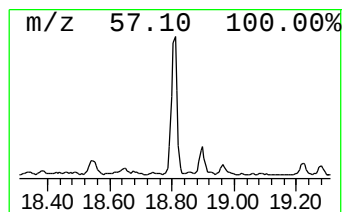
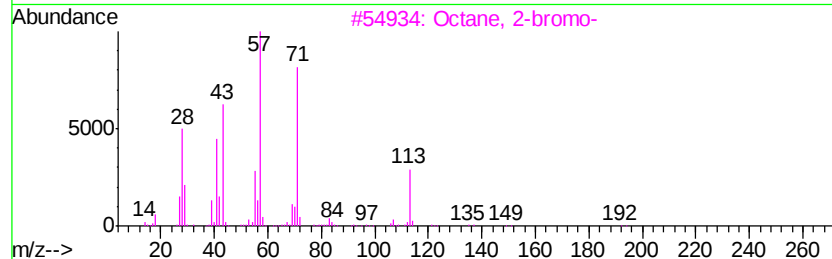
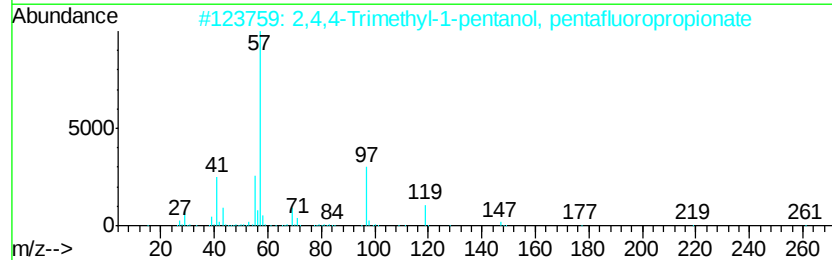
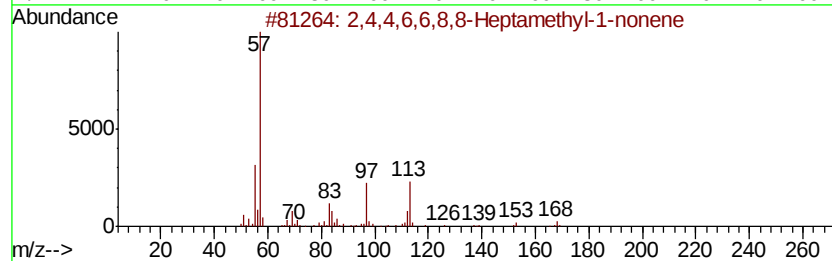
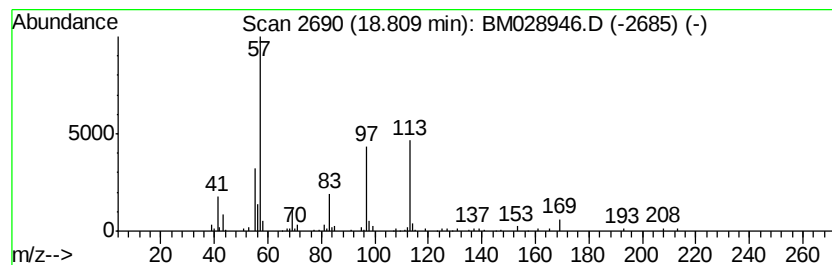
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.81	5.86 ng/ul	61397	Phenanthrene-d10	17.17

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	72
2			2,4,4-Trimethyl-1-pentanol, pent...	276	C11H17F5O2	1000365-51-0	38
3			Octane, 2-bromo-	192	C8H17Br	000557-35-7	37
4			Octane, 2-bromo-	192	C8H17Br	000557-35-7	37
5			1-Pentanol, 3-methyl-2-propyl-	144	C9H20O	054004-40-9	28



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM022721\
Data File : BM028946.D
Acq On : 27 Feb 2021 08:05
Operator : CG/JU
Sample : M1498-02
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
CB8G1

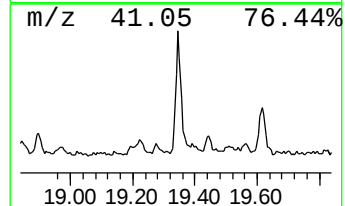
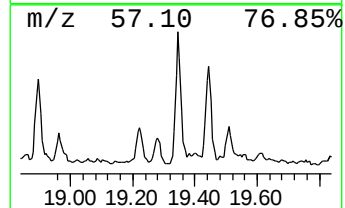
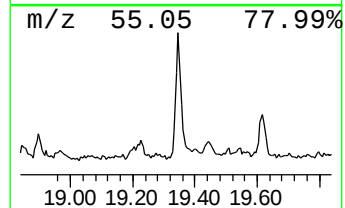
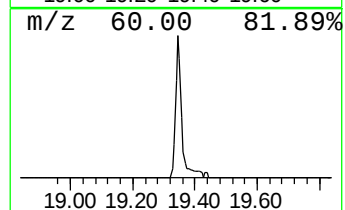
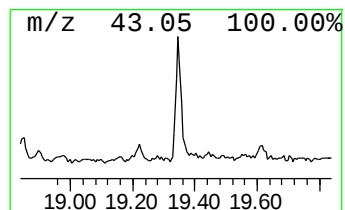
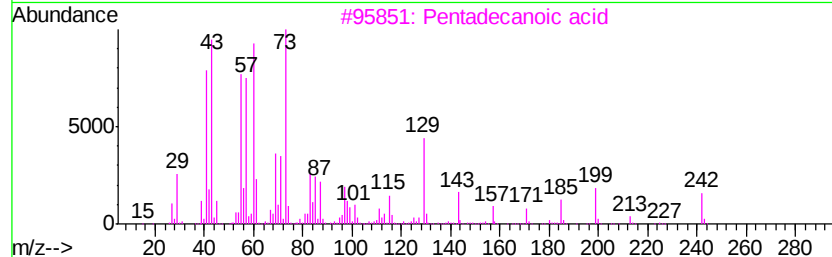
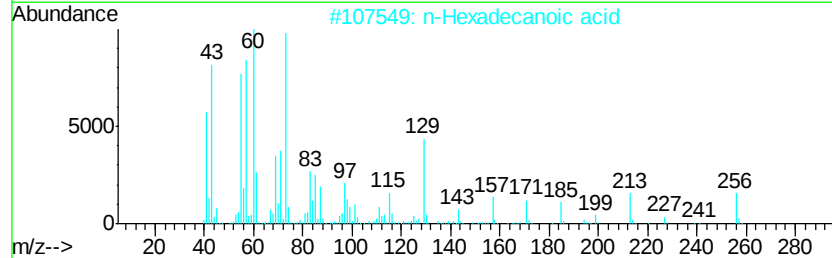
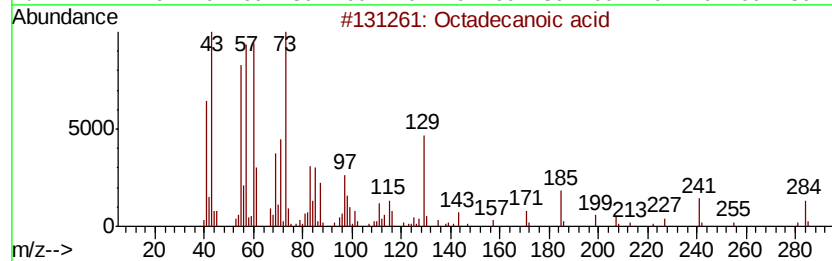
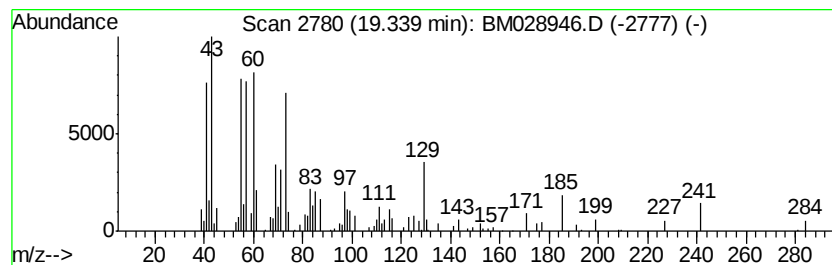
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 Octadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.34	7.11 ng/ul	79369	Chrysene-d12	21.34

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecanoic acid	284	C18H36O2	000057-11-4	95
2		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	81
3		Pentadecanoic acid	242	C15H30O2	001002-84-2	62
4		Tetradecanoic acid	228	C14H28O2	000544-63-8	58
5		Pentadecanoic acid	242	C15H30O2	001002-84-2	46



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM022721\
Data File : BM028946.D
Acq On : 27 Feb 2021 08:05
Operator : CG/JU
Sample : M1498-02
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
CB8G1

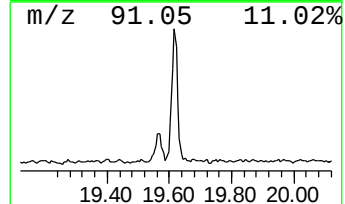
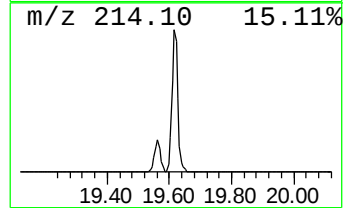
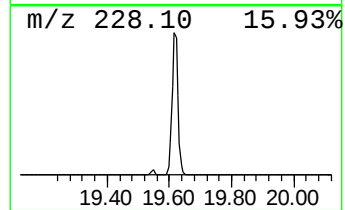
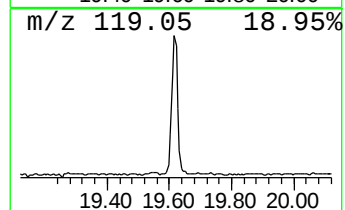
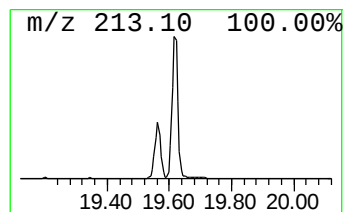
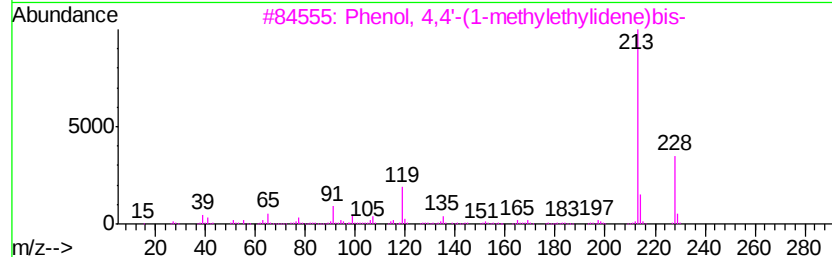
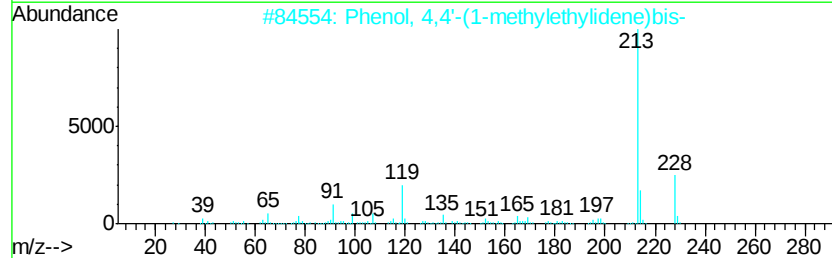
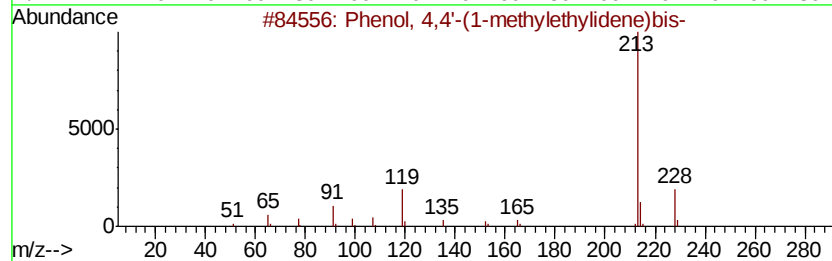
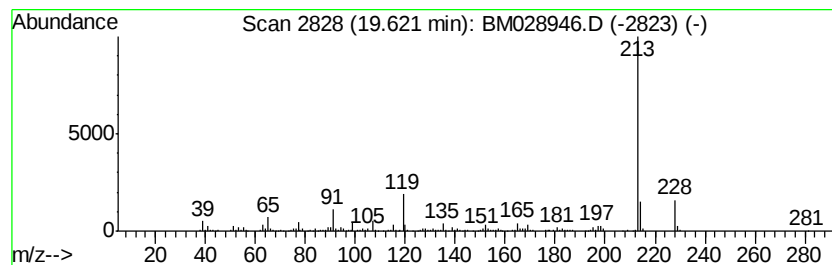
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 Phenol, 4,4'-(1-methylethyl... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.62	18.88 ng/ul	210806	Chrysene-d12	21.34

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenol, 4,4'-(1-methylethylidene)bis-	228	C15H16O2	000080-05-7	97
2		Phenol, 4,4'-(1-methylethylidene)bis-	228	C15H16O2	000080-05-7	95
3		Phenol, 4,4'-(1-methylethylidene)bis-	228	C15H16O2	000080-05-7	90
4		2H,8H-Benzo[1,2-b:3,4-b']dipyran...	228	C14H12O3	000523-59-1	58
5		2H,8H-Benzo[1,2-b:5,4-b']dipyran...	228	C14H12O3	000553-19-5	58



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM022721\
Data File : BM028946.D
Acq On : 27 Feb 2021 08:05
Operator : CG/JU
Sample : M1498-02
Misc :
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Instrument :
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ClientSampleId :
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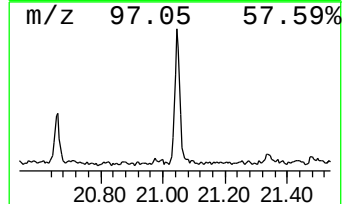
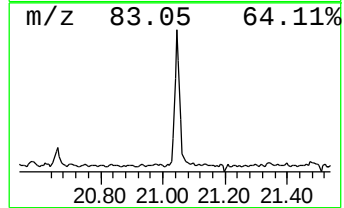
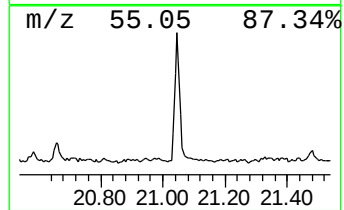
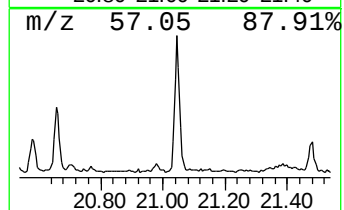
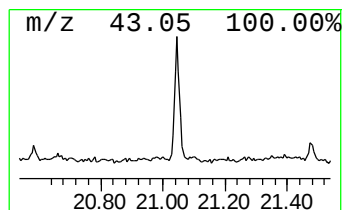
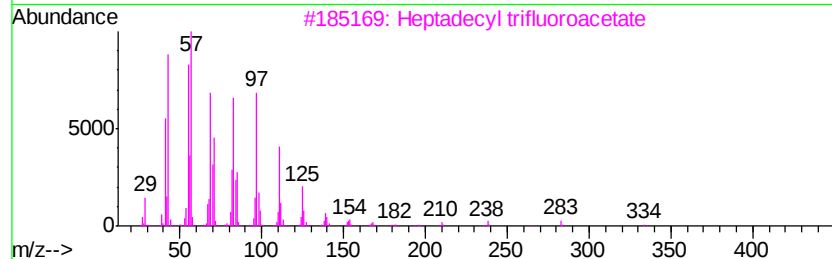
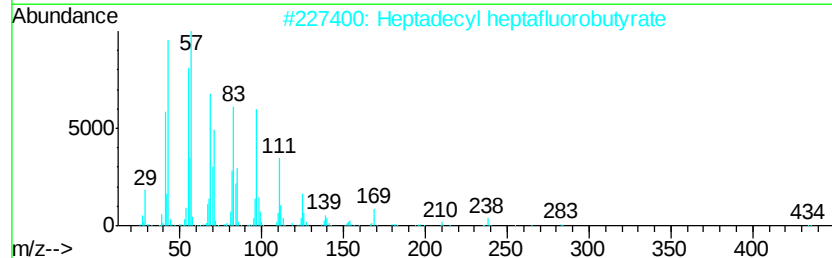
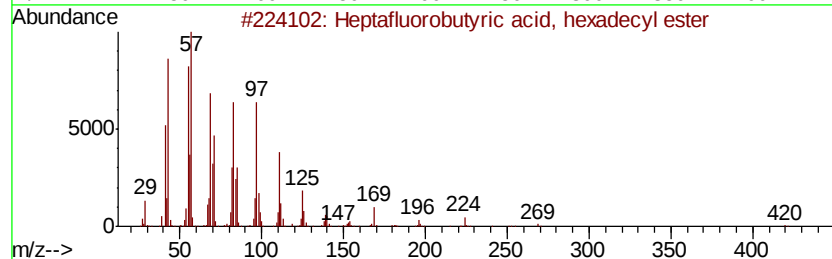
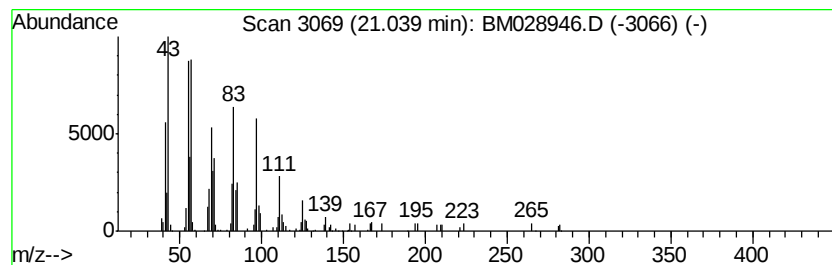
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 Heptafluorobutyric acid, he... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.04	6.83 ng/ul	76285	Chrysene-d12	21.34

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptafluorobutyric acid, hexadec...	438	C20H33F7O2	006385-15-5	95
2		Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	95
3		Heptadecyl trifluoroacetate	352	C19H35F3O2	1000351-87-0	95
4		Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	95
5		Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	95



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM022721\
Data File : BM028946.D
Acq On : 27 Feb 2021 08:05
Operator : CG/JU
Sample : M1498-02
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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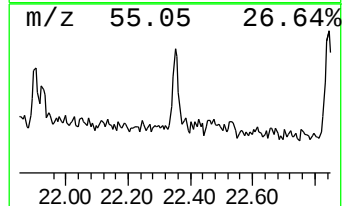
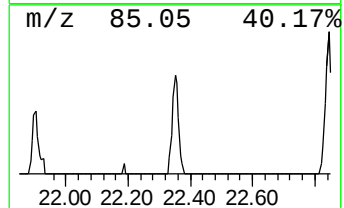
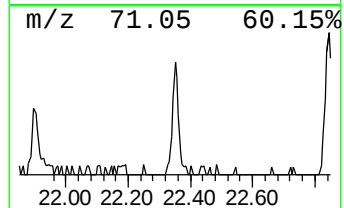
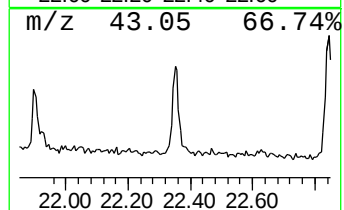
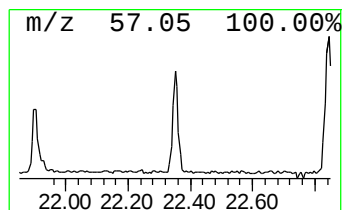
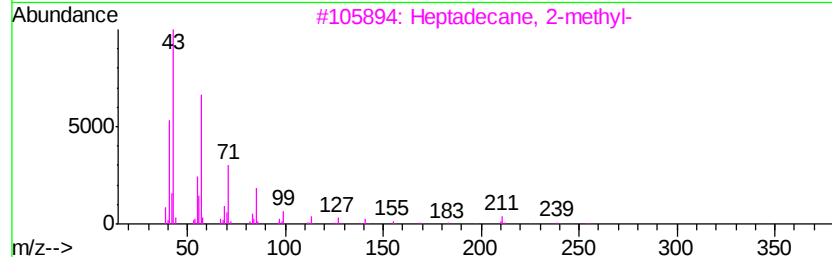
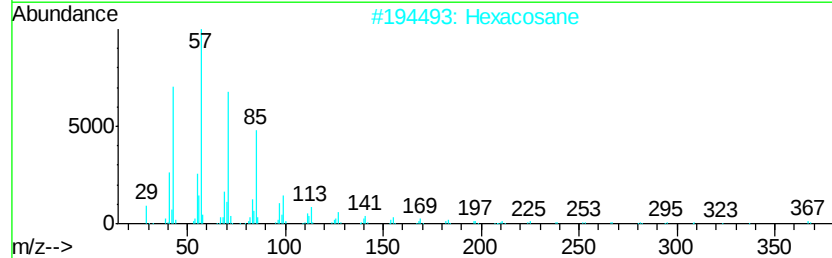
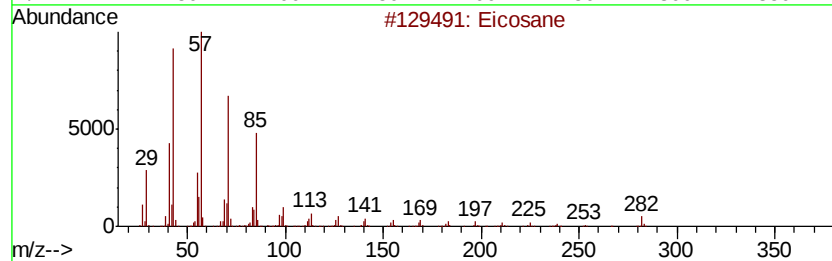
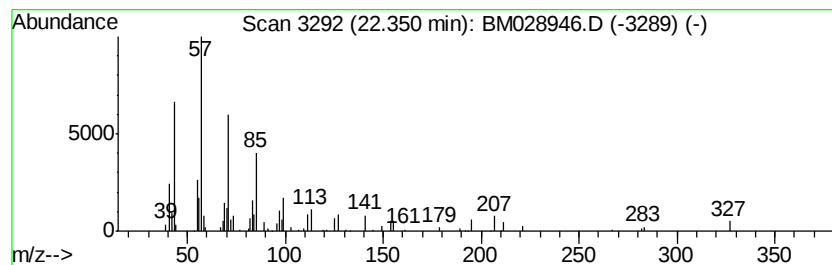
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 (DEL) Alkane: Straight-Chai... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.35	2.31 ng/ul	25746	Chrysene-d12	21.34

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	96
2		Hexacosane	366	C26H54	000630-01-3	87
3		Heptadecane, 2-methyl-	254	C18H38	001560-89-0	87
4		Pentacosane	352	C25H52	000629-99-2	87
5		Hentriacontane	437	C31H64	000630-04-6	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM022721\
Data File : BM028946.D
Acq On : 27 Feb 2021 08:05
Operator : CG/JU
Sample : M1498-02
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
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ClientSampleId :
CB8G1

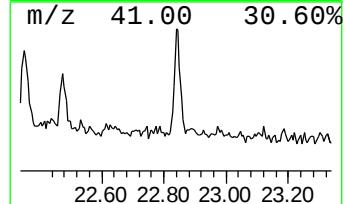
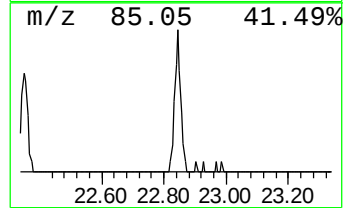
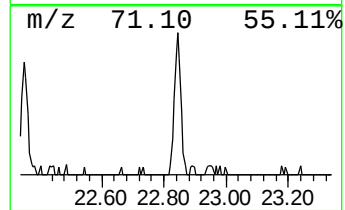
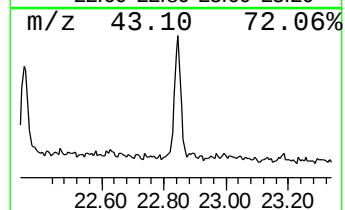
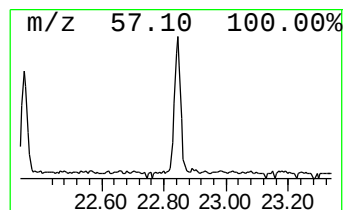
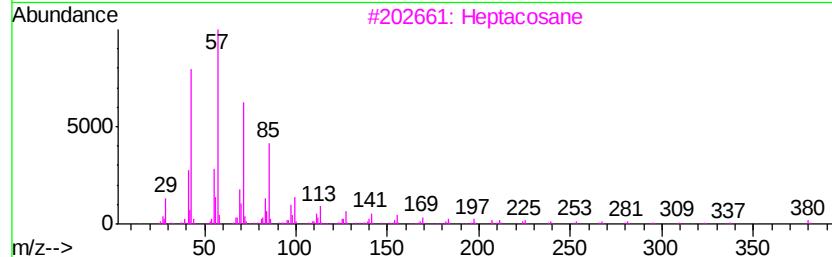
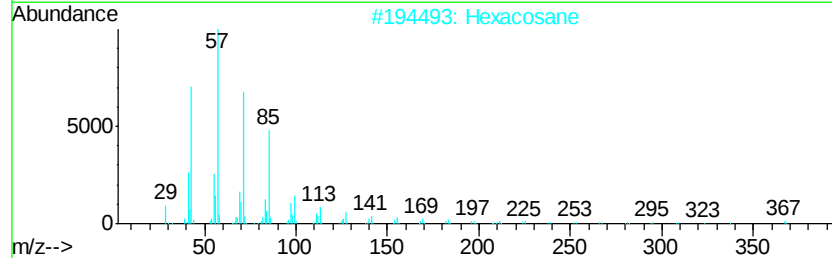
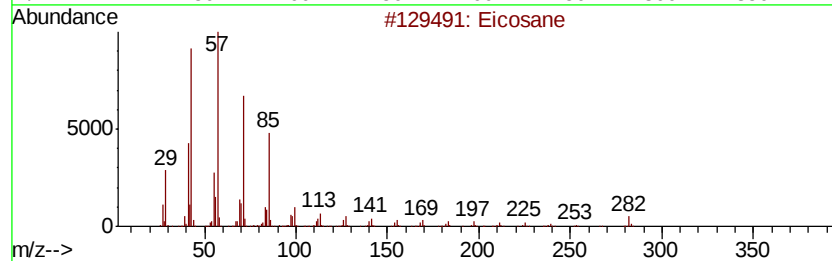
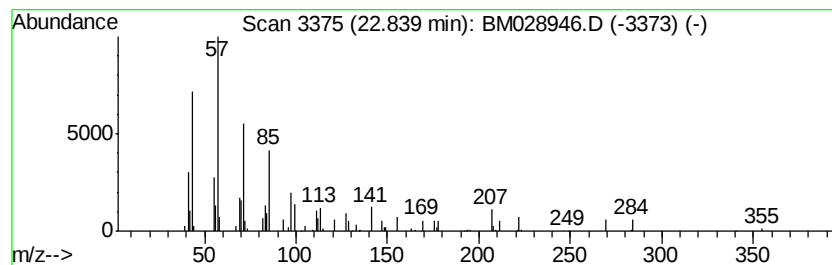
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Quant Title : SVOA CALIBRATION

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

Peak Number 15 (DEL) Alkane: Straight-Chai... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.84	2.71 ng/ul	28269	Perylene-d12	23.54

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	97
2		Hexacosane	366	C26H54	000630-01-3	81
3		Heptacosane	380	C27H56	000593-49-7	81
4		Hexacosane	366	C26H54	000630-01-3	81
5		Hentriacontane	437	C31H64	000630-04-6	81



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM022721\
Data File : BM028946.D
Acq On : 27 Feb 2021 08:05
Operator : CG/JU
Sample : M1498-02
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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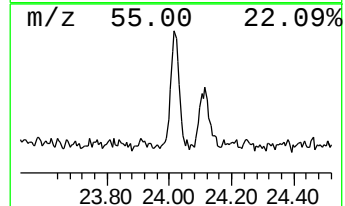
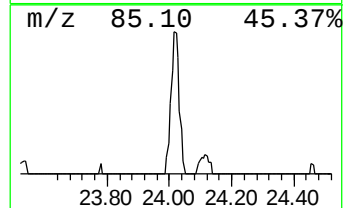
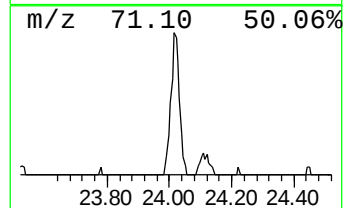
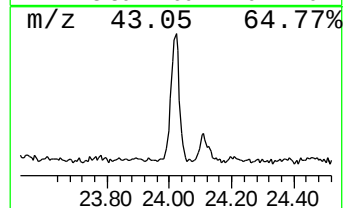
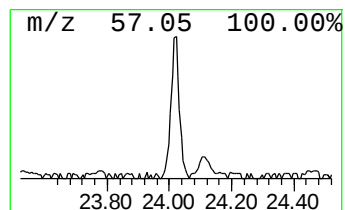
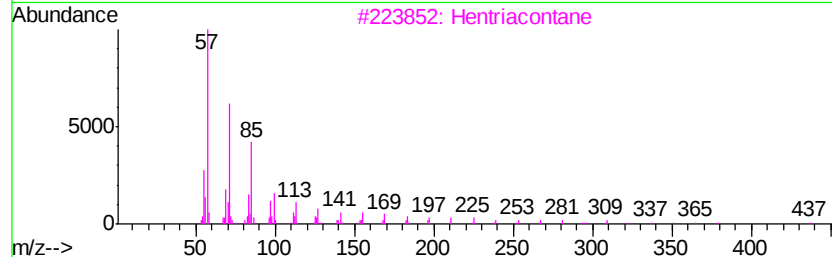
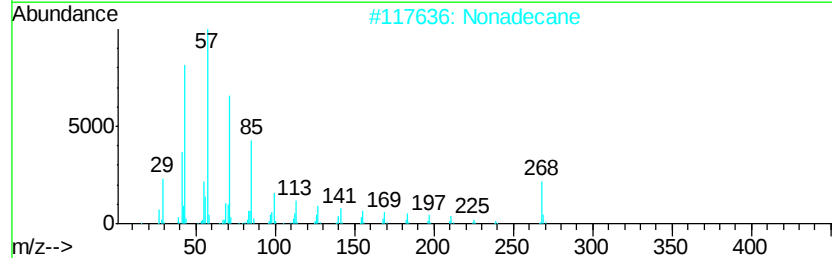
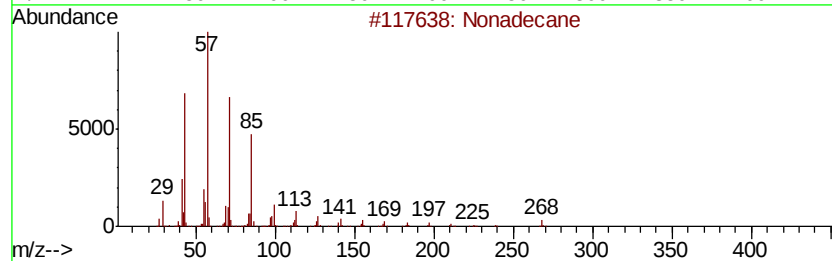
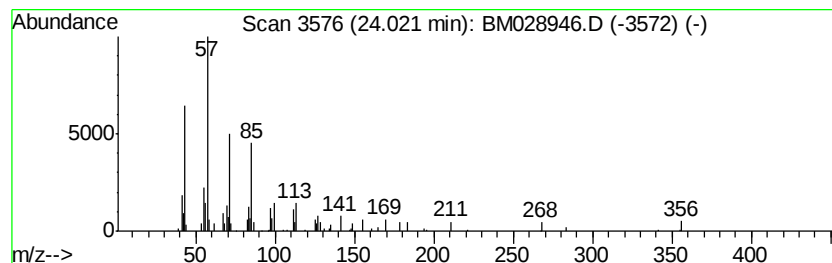
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.02	3.97 ng/ul	41403	Perylene-d12	23.54

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonadecane	268	C19H40	000629-92-5	94
2		Nonadecane	268	C19H40	000629-92-5	94
3		Hentriacontane	437	C31H64	000630-04-6	87
4		Hentriacontane	437	C31H64	000630-04-6	86
5		Docosane, 11-butyl-	366	C26H54	013475-76-8	83



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM022721\
 Data File : BM028946.D
 Acq On : 27 Feb 2021 08:05
 Operator : CG/JU
 Sample : M1498-02
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
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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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
1,3-Oxathiolane	4.35	5.7	ng/ul	30905	1	7.77	109364	20.0
Benzene, chloro-	5.04	7.0	ng/ul	38068	1	7.77	109364	20.0
Benzenamine, 3-me...	8.77	3.0	ng/ul	16638	1	7.77	109364	20.0
Phenol, p-tert-bu...	11.93	2.4	ng/ul	18434	2	10.57	155769	20.0
unknown-01	14.36	2.5	ng/ul	24805	3	14.43	195793	20.0
Diethyltoluamide	15.19	8.5	ng/ul	83110	3	14.43	195793	20.0
Benzenesulfonamid...	15.96	12.3	ng/ul	129068	4	17.17	209449	20.0
Benzenesulfonamid...	16.47	5.9	ng/ul	62008	4	17.17	209449	20.0
n-Hexadecanoic acid	18.07	12.6	ng/ul	132082	4	17.17	209449	20.0
(DEL) Alkane: Str...	18.81	5.9	ng/ul	61397	4	17.17	209449	20.0
Octadecanoic acid	19.34	7.1	ng/ul	79369	5	21.34	223334	20.0
Phenol, 4,4'-(1-m...	19.62	18.9	ng/ul	210806	5	21.34	223334	20.0
Heptafluorobutyri...	21.04	6.8	ng/ul	76285	5	21.34	223334	20.0
(DEL) Alkane: Str...	22.35	2.3	ng/ul	25746	5	21.34	223334	20.0
(DEL) Alkane: Str...	22.84	2.7	ng/ul	28269	6	23.54	208717	20.0
(DEL) Alkane: Str...	24.02	4.0	ng/ul	41403	6	23.54	208717	20.0