

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM070220\
 Data File : BM026651.D
 Acq On : 02 Jul 2020 13:08
 Operator : JU/CG
 Sample : L3131-01
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BG2L6

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-BM063020MA.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	2.951	8	11	18	rVB	31976	40020	3.03%	0.128%
2	3.481	98	101	105	rVV3	11778	13622	1.03%	0.044%
3	3.557	109	114	118	rBV2	11203	17249	1.31%	0.055%
4	3.640	125	128	132	rVB2	8550	11792	0.89%	0.038%
5	3.869	163	167	171	rBV3	11120	13353	1.01%	0.043%
6	4.534	276	280	288	rVV	18561	25126	1.90%	0.080%
7	6.057	533	539	544	rBV	73385	101940	7.72%	0.326%
8	6.351	583	589	596	rBV7	8069	13731	1.04%	0.044%
9	6.428	596	602	612	rBV	34084	77673	5.88%	0.248%
10	6.557	618	624	637	rVB	399909	645295	48.85%	2.063%
11	6.710	644	650	661	rBV	325060	492454	37.28%	1.575%
12	6.787	661	663	669	rVB5	7756	11751	0.89%	0.038%
13	6.892	675	681	696	rBV	419352	638249	48.32%	2.041%
14	7.351	752	759	767	rBV	438005	644716	48.81%	2.061%
15	7.575	793	797	802	rBV3	8958	13710	1.04%	0.044%
16	8.075	876	882	894	rVV	396318	640762	48.51%	2.049%
17	8.181	894	900	907	rVV	11955	24488	1.85%	0.078%
18	8.498	946	954	965	rBV	383930	608866	46.09%	1.947%
19	9.210	1068	1075	1086	rBV	308009	505757	38.29%	1.617%
20	9.751	1160	1167	1182	rBV	389007	698467	52.88%	2.233%
21	10.104	1220	1227	1239	rBV	519572	851615	64.47%	2.723%
22	10.263	1247	1254	1270	rBV	288465	590538	44.71%	1.888%
23	12.198	1578	1583	1598	rBV2	35176	77526	5.87%	0.248%
24	12.351	1601	1609	1615	rVB5	8187	14852	1.12%	0.047%
25	12.616	1650	1654	1661	rVB5	6585	10155	0.77%	0.032%
26	12.880	1689	1699	1702	rBV5	25171	56504	4.28%	0.181%
27	12.927	1702	1707	1714	rVB	206155	311392	23.57%	0.996%
28	13.327	1770	1775	1779	rVV3	16728	24765	1.87%	0.079%
29	13.439	1784	1794	1798	rVV	704785	961486	72.79%	3.074%
30	13.486	1798	1802	1812	rVV	195390	274286	20.77%	0.877%
31	13.698	1831	1838	1848	rBV	700391	1033887	78.27%	3.306%
32	13.980	1875	1886	1888	rBV	283699	416514	31.53%	1.332%
33	14.010	1888	1891	1899	rVB	715312	963290	72.93%	3.080%
34	14.104	1903	1907	1914	rVB6	14612	27899	2.11%	0.089%

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Integrator: RTE
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Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM063020MA.M
 Title : SVOA CALIBRATION

35	14.268	1928	1935	1947	rBV	158442	405466	30.70%	1.296%
36	14.480	1970	1971	1980	rVB8	5744	10509	0.80%	0.034%
37	14.663	1999	2002	2006	rVV6	9330	11002	0.83%	0.035%
38	14.757	2014	2018	2025	rBV2	47187	70714	5.35%	0.226%
39	14.851	2030	2034	2040	rVV2	8422	13550	1.03%	0.043%
40	14.910	2040	2044	2049	rVB2	10379	12300	0.93%	0.039%
41	15.015	2056	2062	2072	rBV	863764	1178751	89.24%	3.769%
42	15.098	2073	2076	2082	rVV6	7640	11757	0.89%	0.038%
43	15.162	2082	2087	2100	rVV3	274173	540001	40.88%	1.727%
44	15.257	2100	2103	2107	rVV2	12193	17831	1.35%	0.057%
45	15.315	2109	2113	2117	rVV3	8145	12980	0.98%	0.042%
46	15.362	2117	2121	2125	rVV5	7364	10498	0.79%	0.034%
47	15.515	2143	2147	2153	rVB2	26452	34577	2.62%	0.111%
48	15.610	2157	2163	2166	rBV7	6009	11608	0.88%	0.037%
49	15.715	2178	2181	2187	rVB7	4937	10053	0.76%	0.032%
50	15.780	2187	2192	2193	rBV2	14035	18420	1.39%	0.059%
51	15.804	2193	2196	2205	rVB4	18166	31226	2.36%	0.100%
52	15.880	2206	2209	2215	rBV7	8456	14120	1.07%	0.045%
53	16.021	2226	2233	2240	rVB	28839	44394	3.36%	0.142%
54	16.186	2257	2261	2264	rBV4	11016	14017	1.06%	0.045%
55	16.433	2297	2303	2314	rBV	517561	780504	59.09%	2.496%
56	16.568	2321	2326	2331	rVV4	21330	32781	2.48%	0.105%
57	16.621	2331	2335	2342	rVB	21714	37105	2.81%	0.119%
58	16.768	2354	2360	2365	rBV	827163	1110746	84.09%	3.552%
59	16.809	2365	2367	2371	rVB	31344	34711	2.63%	0.111%
60	16.868	2371	2377	2386	rBV	855950	1206776	91.36%	3.859%
61	17.109	2414	2418	2422	rBV4	19006	28173	2.13%	0.090%
62	17.151	2422	2425	2429	rVV4	14017	18938	1.43%	0.061%
63	17.304	2449	2451	2455	rVB	23449	26351	1.99%	0.084%
64	17.398	2463	2467	2472	rVB8	10253	15552	1.18%	0.050%
65	17.709	2514	2520	2526	rBV	205655	309720	23.45%	0.990%
66	17.768	2526	2530	2534	rVB	40819	66224	5.01%	0.212%
67	17.998	2565	2569	2577	rVB	52860	68033	5.15%	0.218%
68	18.186	2595	2601	2605	rBV2	24470	46937	3.55%	0.150%
69	18.286	2616	2618	2622	rVB5	18918	19874	1.50%	0.064%
70	18.456	2642	2647	2653	rBV	172129	236742	17.92%	0.757%
71	18.545	2659	2662	2664	rBV	27491	29873	2.26%	0.096%

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72	18.574	2664	2667	2672	rVV2	62066	97124	7.35%	0.311%
73	18.645	2675	2679	2687	rVB	106625	130319	9.87%	0.417%
74	18.809	2704	2707	2709	rBV3	20001	22191	1.68%	0.071%
75	18.839	2709	2712	2715	rVV	48283	61829	4.68%	0.198%
76	18.868	2715	2717	2720	rVV2	32247	45425	3.44%	0.145%
77	18.898	2720	2722	2728	rVB	61987	79064	5.99%	0.253%
78	19.009	2736	2741	2747	rBV2	97354	171941	13.02%	0.550%
79	19.174	2764	2769	2776	rBV2	984961	1320898	100.00%	4.223%
80	19.233	2776	2779	2783	rVB	193710	201494	15.25%	0.644%
81	19.639	2845	2848	2852	rBV2	30075	42800	3.24%	0.137%
82	19.745	2863	2866	2868	rBV3	33031	42706	3.23%	0.137%
83	19.774	2868	2871	2875	rVB	292819	307089	23.25%	0.982%
84	20.209	2943	2945	2950	rVB2	39891	45177	3.42%	0.144%
85	20.274	2952	2956	2960	rVV	380659	434928	32.93%	1.391%
86	20.339	2964	2967	2973	rVB	114070	135938	10.29%	0.435%
87	20.556	3000	3004	3012	rBV	805265	961883	72.82%	3.076%
88	20.739	3031	3035	3046	rVB	914672	1180641	89.38%	3.775%
89	20.939	3066	3069	3072	rBV2	56299	76534	5.79%	0.245%
90	20.986	3072	3077	3081	rVV	953352	1166177	88.29%	3.729%
91	21.180	3106	3110	3114	rBV	532654	572688	43.36%	1.831%
92	21.597	3177	3181	3189	rBV	557890	716678	54.26%	2.292%
93	22.021	3249	3253	3258	rBV	546688	664231	50.29%	2.124%
94	22.486	3328	3332	3337	rVB	579123	730581	55.31%	2.336%
95	22.938	3403	3409	3414	rBV	687943	1143206	86.55%	3.655%
96	22.997	3415	3419	3424	rVV	452461	666332	50.45%	2.131%
97	23.062	3425	3430	3438	rVB2	719299	1238508	93.76%	3.960%
98	23.574	3513	3517	3525	rVB	417574	705649	53.42%	2.256%
99	24.238	3625	3630	3636	rBV	194022	383487	29.03%	1.226%
100	24.480	3667	3671	3682	rVB4	244545	533052	40.36%	1.704%

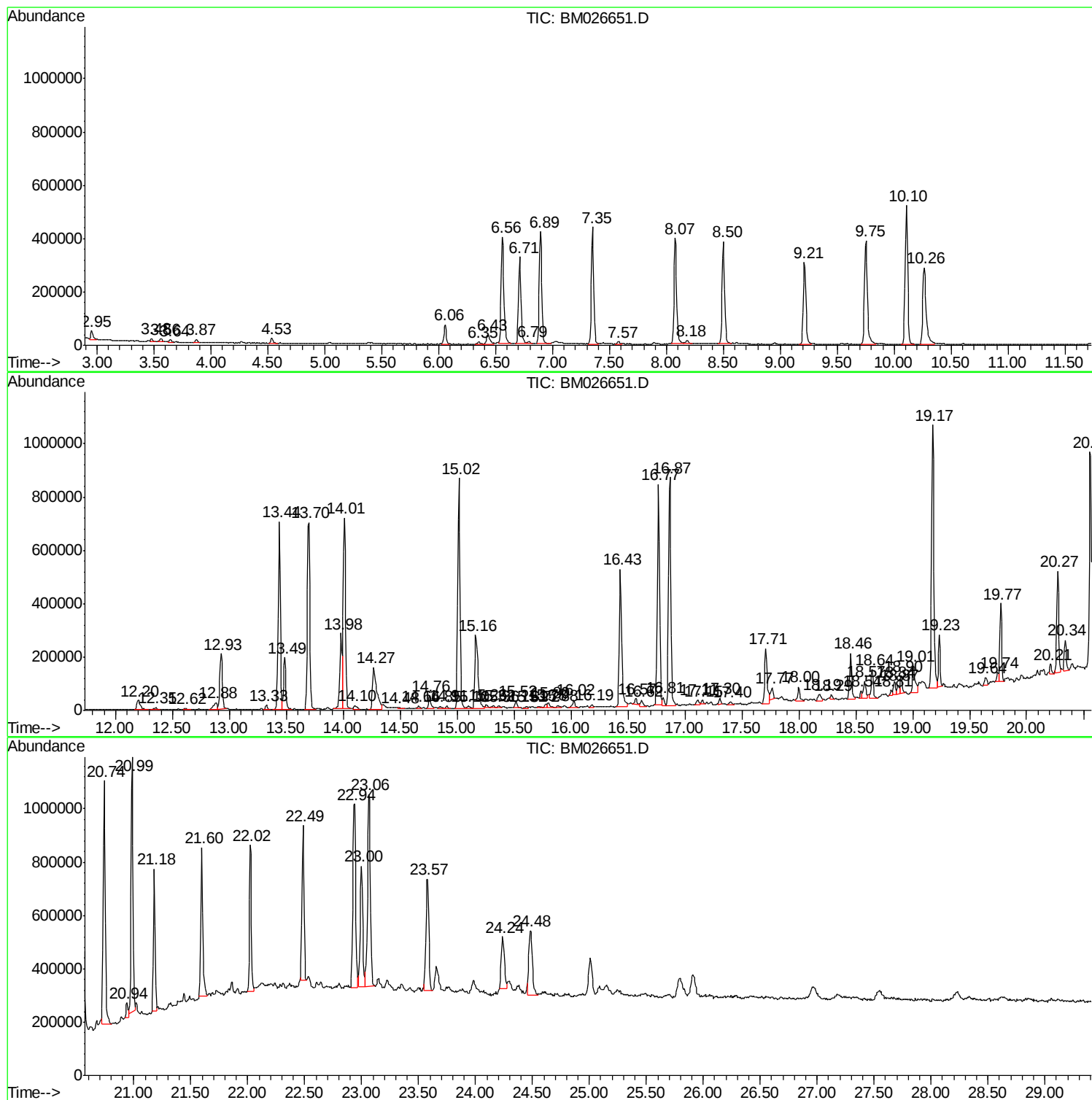
Sum of corrected areas: 31275114

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM063020MA.M
Quant Title : SVOA CALIBRATION

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



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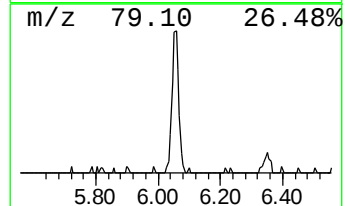
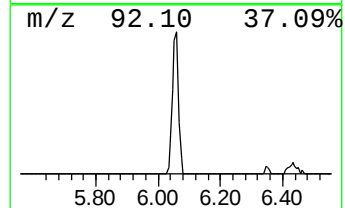
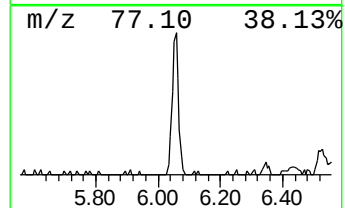
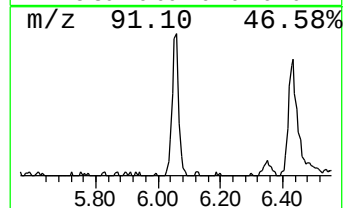
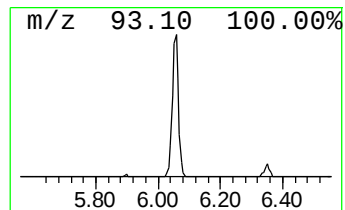
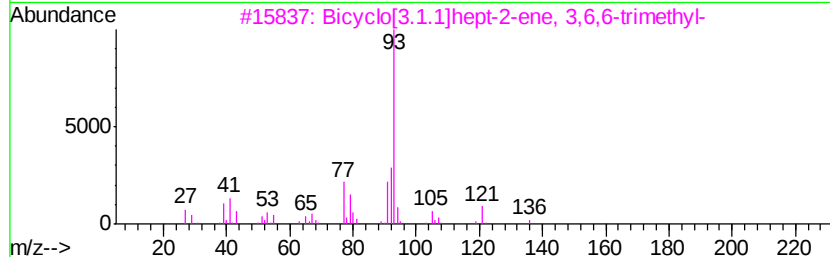
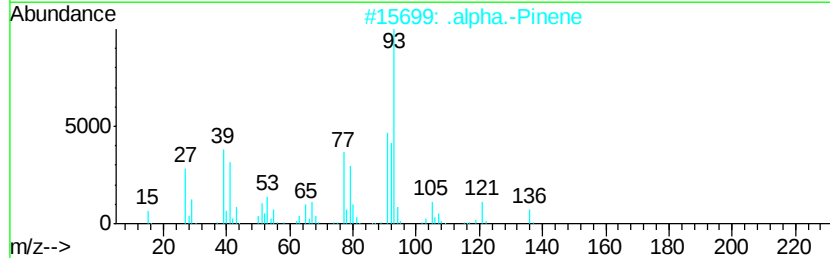
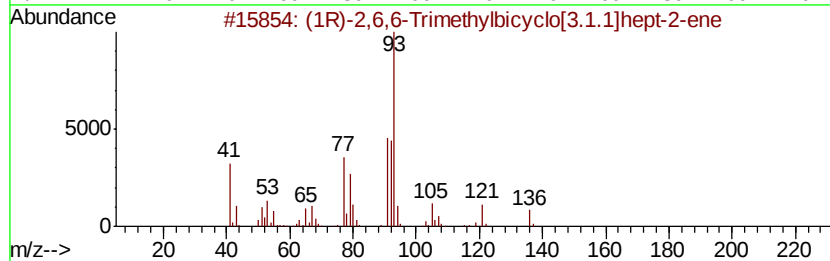
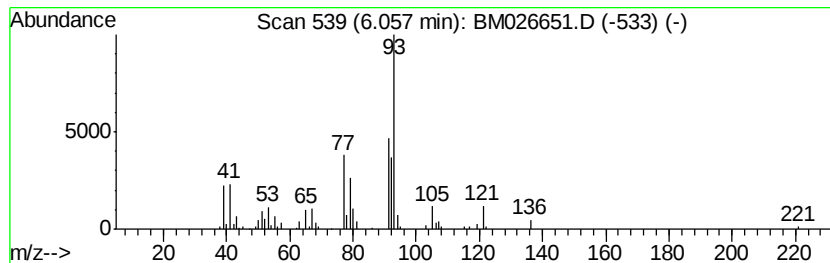
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM063020MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 (1R)-2,6,6-Trimethylbicyclo... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.06	3.16 ng/ul	101940	1,4-Dichlorobenzene-d4	7.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(1R)-2,6,6-Trimethylbicyclo[3.1.1]hept-2-ene	136	C10H16	007785-70-8	96
2		.alpha.-Pinene	136	C10H16	000080-56-8	94
3		Bicyclo[3.1.1]hept-2-ene, 3,6,6-trimethyl-	136	C10H16	004889-83-2	91
4		(1R)-2,6,6-Trimethylbicyclo[3.1.1]hept-2-ene	136	C10H16	007785-70-8	81
5		(1S)-2,6,6-Trimethylbicyclo[3.1.1]hept-2-ene	136	C10H16	007785-26-4	81



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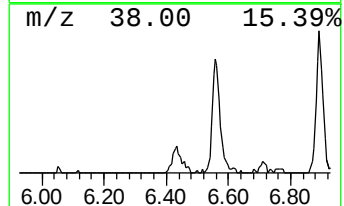
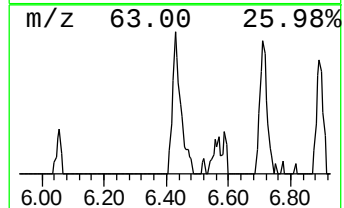
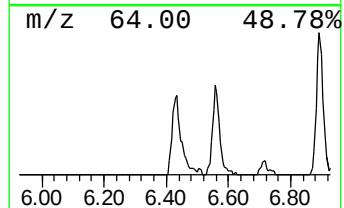
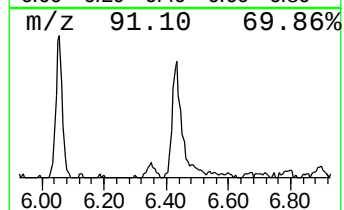
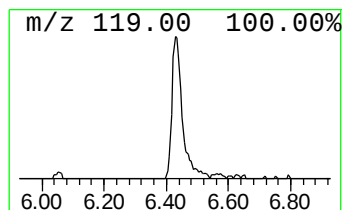
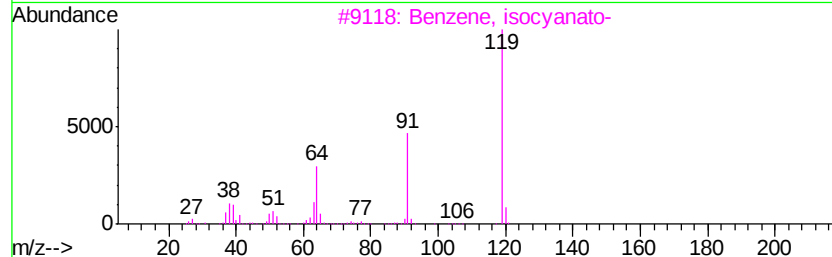
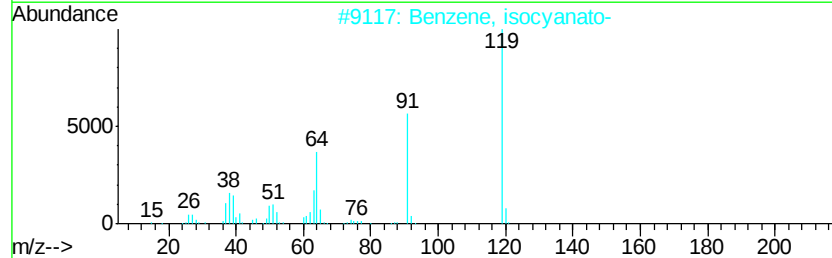
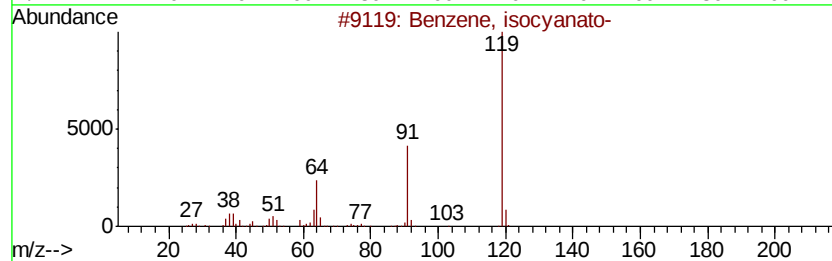
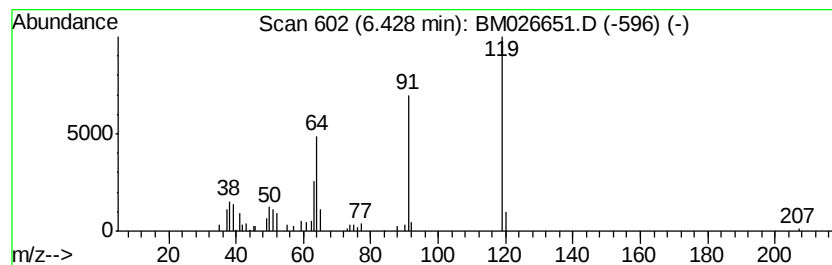
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM063020MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzene, isocyanato- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.43	2.41 ng/ul	77673	1,4-Dichlorobenzene-d4	7.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, isocvanato-	119	C7H5NO	000103-71-9	94
2		Benzene, isocvanato-	119	C7H5NO	000103-71-9	94
3		Benzene, isocvanato-	119	C7H5NO	000103-71-9	94
4		2,1-Benzisoxazole	119	C7H5NO	000271-58-9	91
5		2,1-Benzisoxazole	119	C7H5NO	000271-58-9	90



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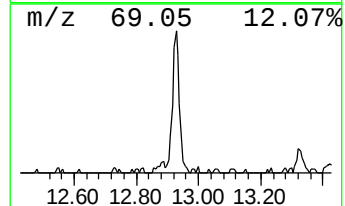
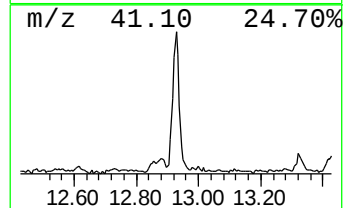
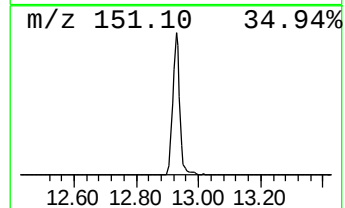
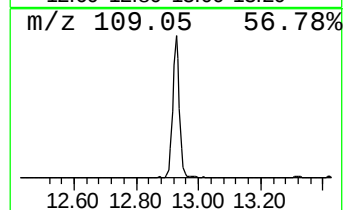
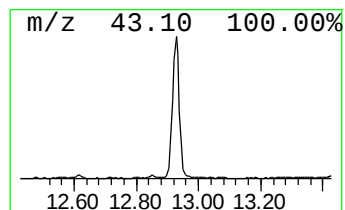
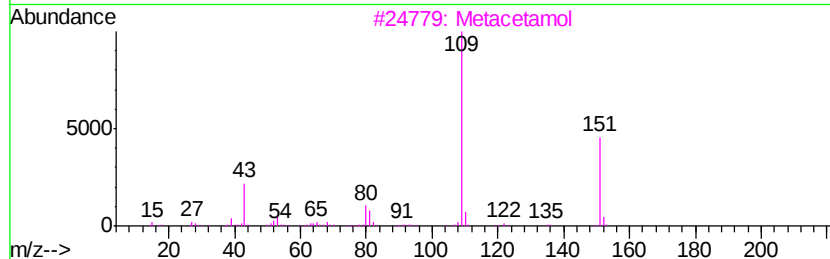
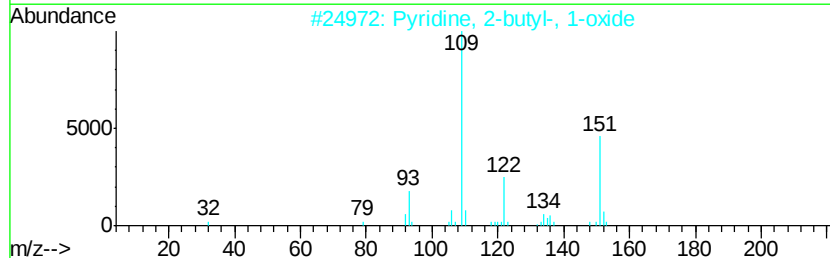
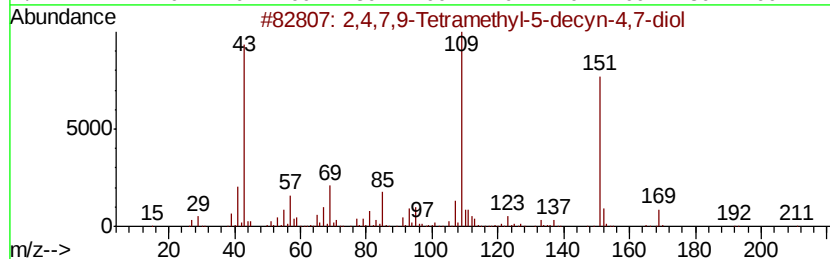
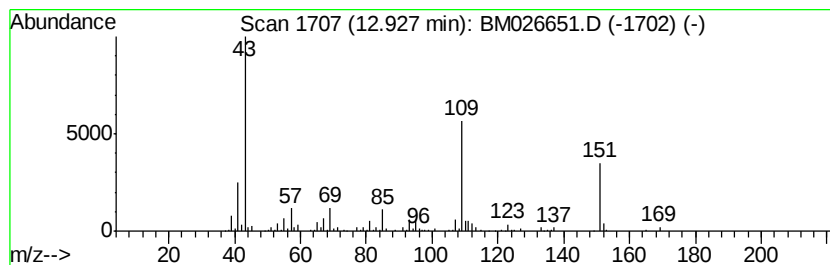
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 2,4,7,9-Tetramethyl-5-decyn... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.93	6.47 ng/ul	311392	Acenaphthene-d10	14.01

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	Pyridine, 2-butyl-, 1-oxide	151	C9H13NO	031396-32-4	50		
3	Metacetamol	151	C8H9NO2	000621-42-1	50		
4	p-Isopropoxyvaniline	151	C9H13NO	007664-66-6	40		
5	1-(1,2,3-Trimethyl-cyclopent-2-e...	152	C10H16O	070987-81-4	38		



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Data File : BM026651.D
Acq On : 02 Jul 2020 13:08
Operator : JU/CG
Sample : L3131-01
Misc :
ALS Vial : 6 Sample Multiplier: 1

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ClientSampleId :
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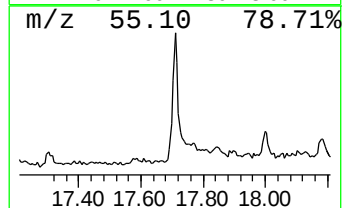
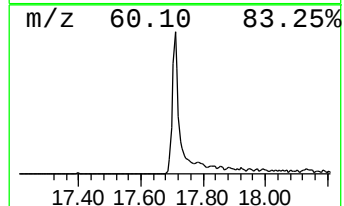
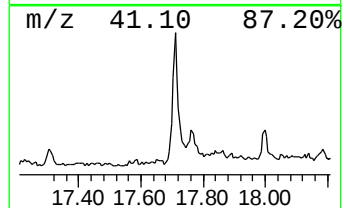
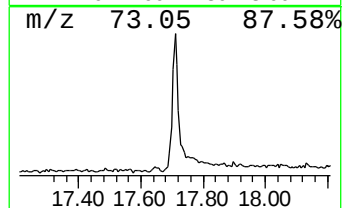
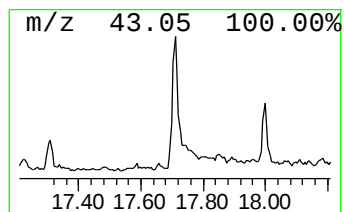
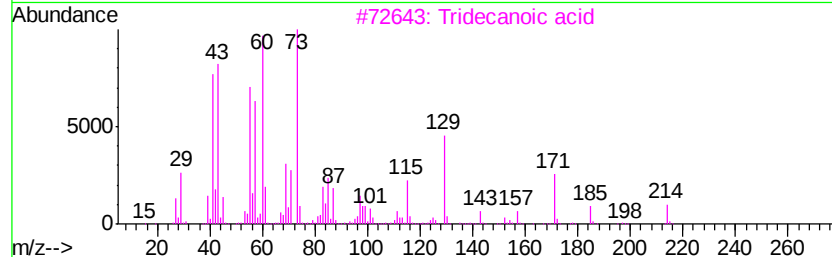
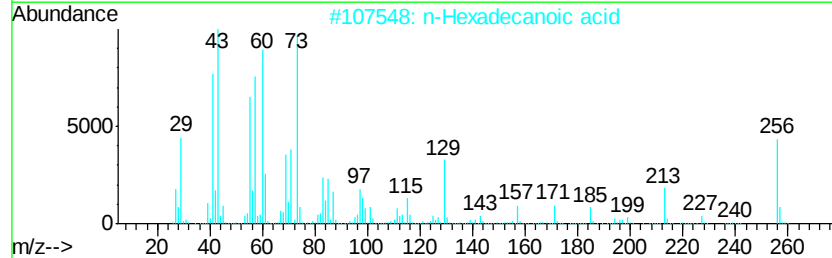
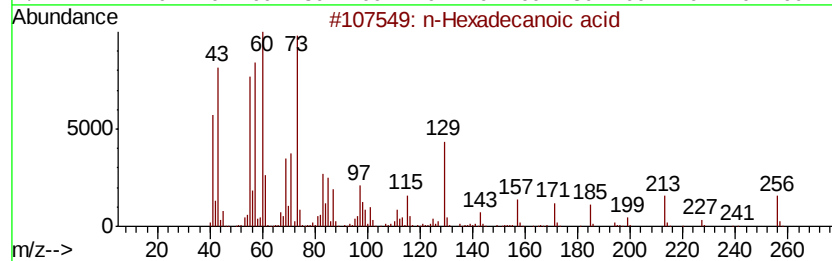
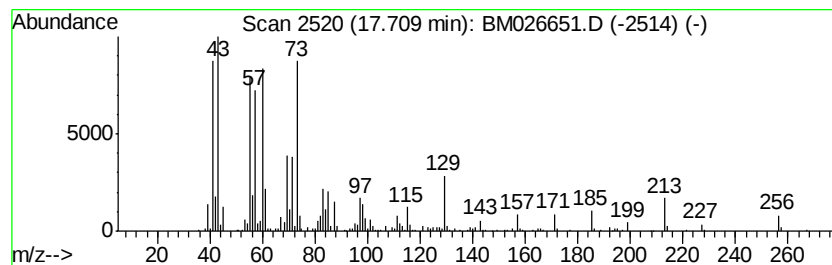
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 n-Hexadecanoic acid Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.71	5.58 ng/ul	309720	Phenanthrene-d10	16.77

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
3		Tridecanoic acid	214	C13H26O2	000638-53-9	94
4		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	91
5		Tetradecanoic acid	228	C14H28O2	000544-63-8	91



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Operator : JU/CG
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ClientSampleId :
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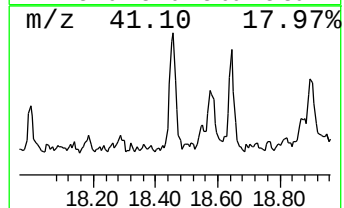
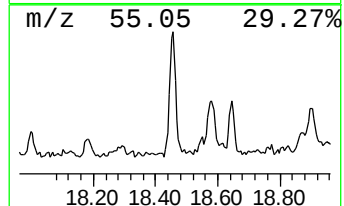
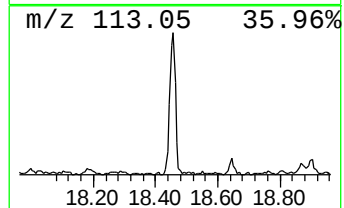
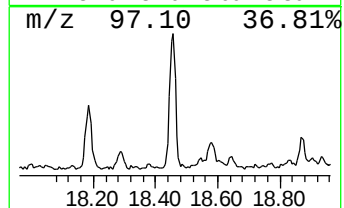
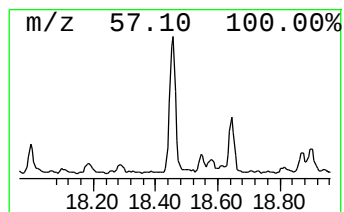
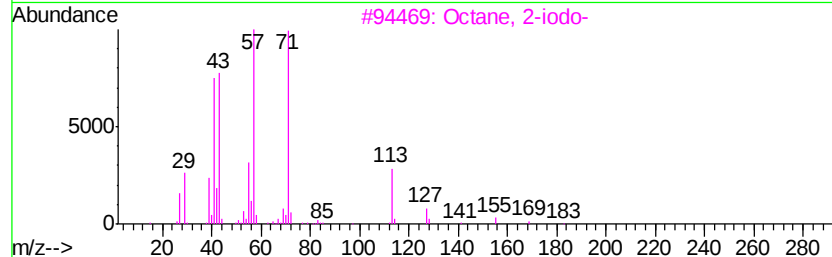
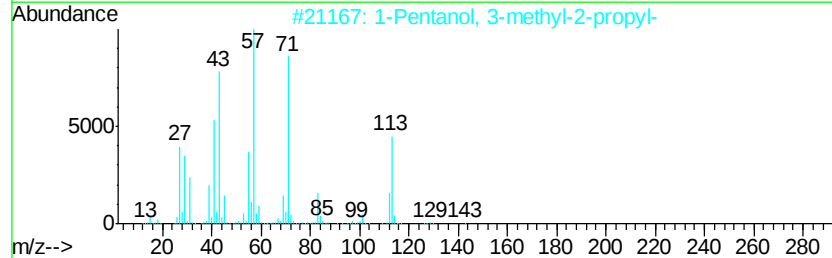
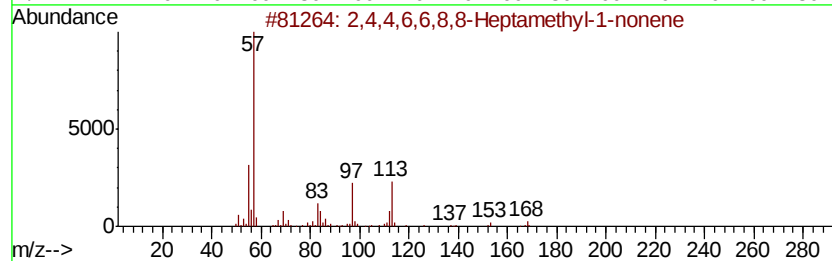
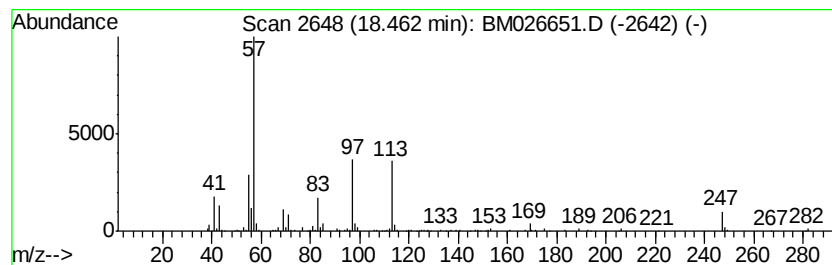
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.46	4.26 ng/ul	236742	Phenanthrene-d10	16.77

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	45		
2	1-Pentanol, 3-methyl-2-propyl-	144	C9H20O	054004-40-9	42		
3	Octane, 2-iodo-	240	C8H17I	000557-36-8	38		
4	2,4,4-Trimethyl-1-pentanol, hept...	326	C12H17F7O2	1000365-01-4	38		
5	2,4,4-Trimethyl-1-pentanol, pent...	276	C11H17F5O2	1000365-51-0	37		



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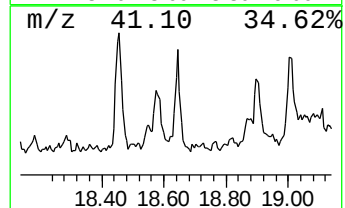
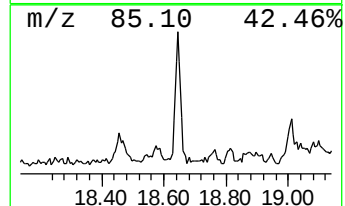
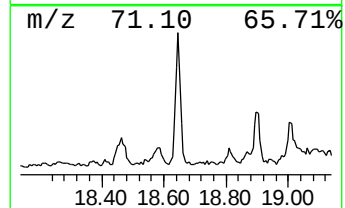
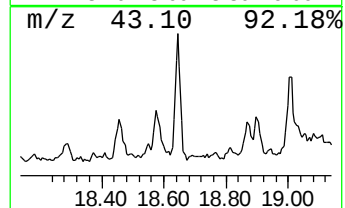
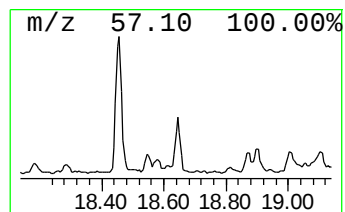
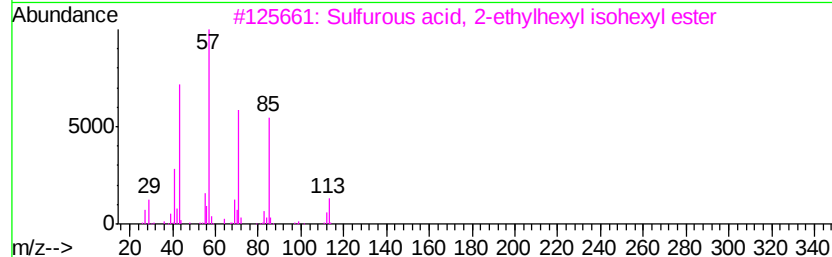
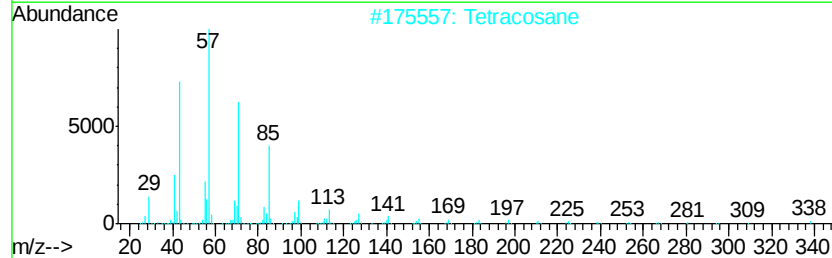
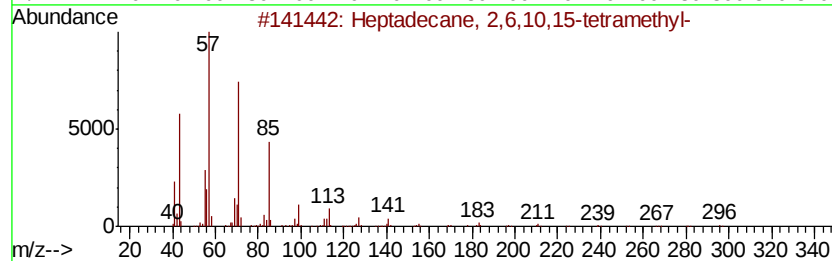
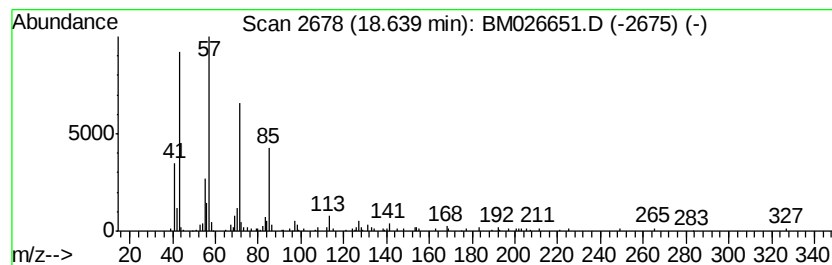
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.64	2.35 ng/ul	130319	Phenanthrene-d10	16.77

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	91
2			Tetracosane	338	C24H50	000646-31-1	90
3			Sulfurous acid, 2-ethylhexyl iso...	278	C14H30O3S	1000309-19-0	80
4			Octadecane, 1-iodo-	380	C18H37I	000629-93-6	80
5			Sulfurous acid, 2-ethylhexyl hex...	278	C14H30O3S	1000309-20-2	78



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM070220\
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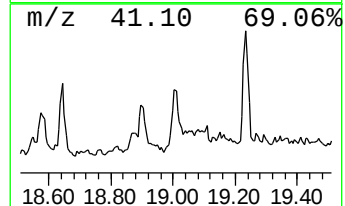
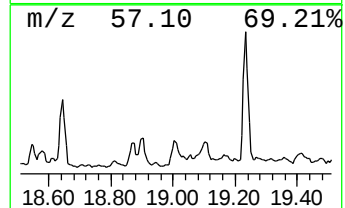
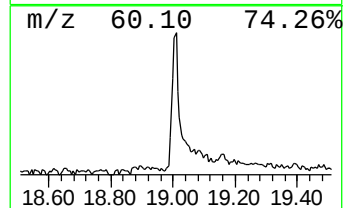
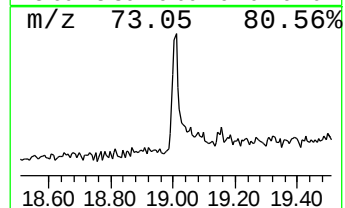
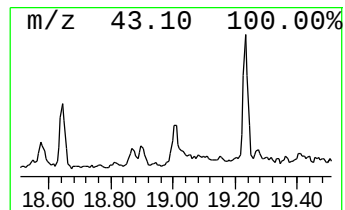
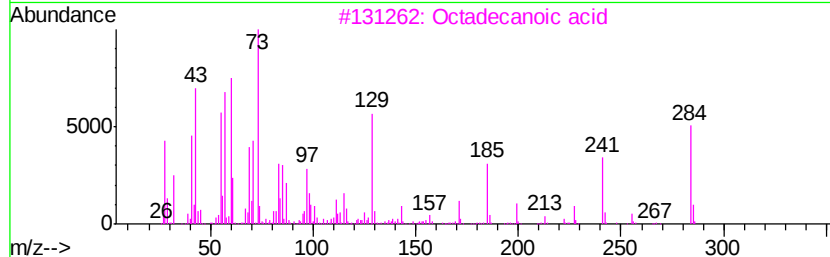
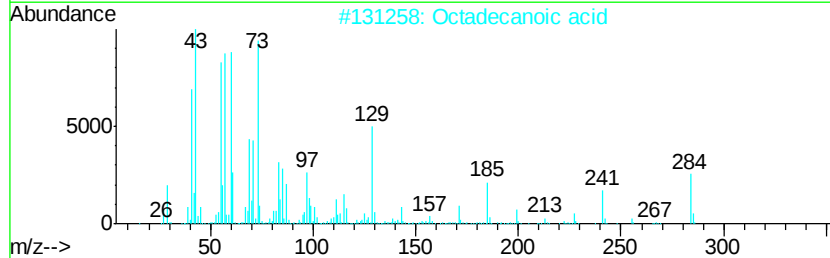
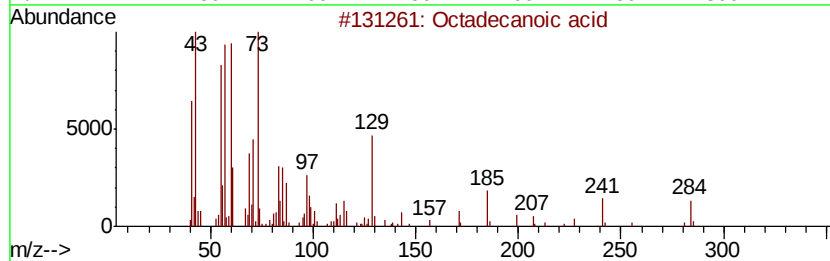
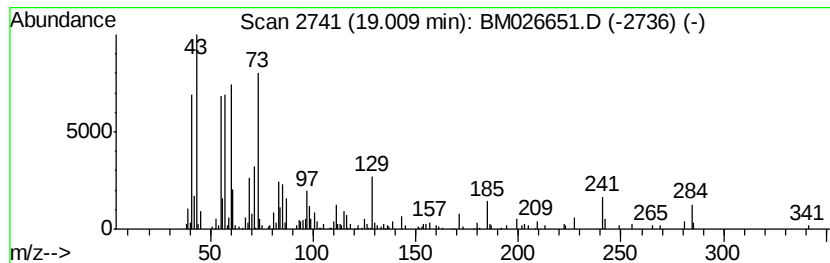
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Octadecanoic acid Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.01	2.95 ng/ul	171941	Chrysene-d12	20.99

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2		Octadecanoic acid	284	C18H36O2	000057-11-4	99
3		Octadecanoic acid	284	C18H36O2	000057-11-4	99
4		Octadecanoic acid	284	C18H36O2	000057-11-4	99
5		Octadecanoic acid	284	C18H36O2	000057-11-4	96



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Data File : BM026651.D
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Misc :
ALS Vial : 6 Sample Multiplier: 1

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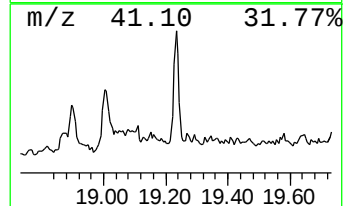
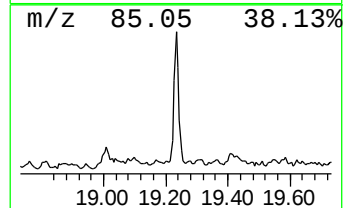
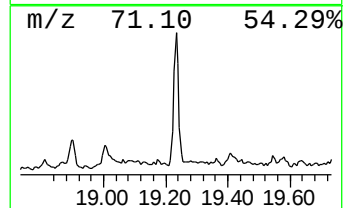
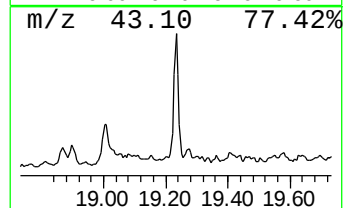
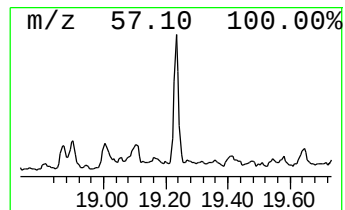
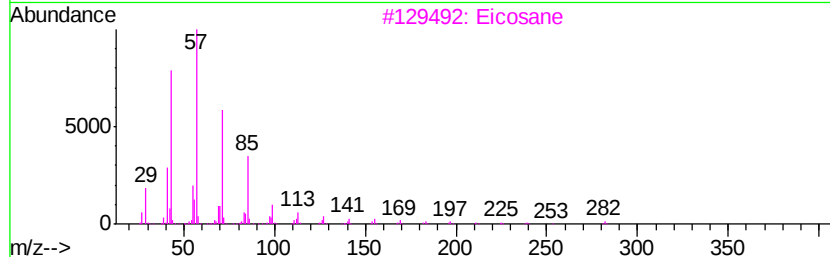
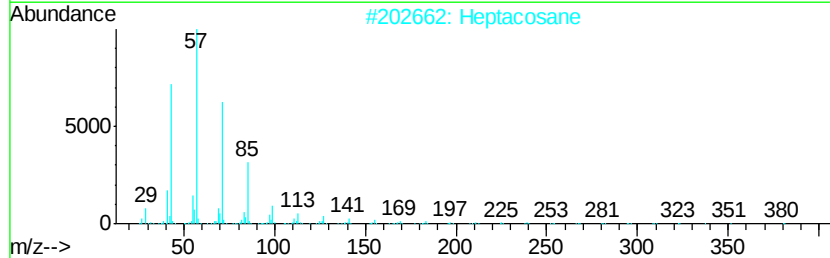
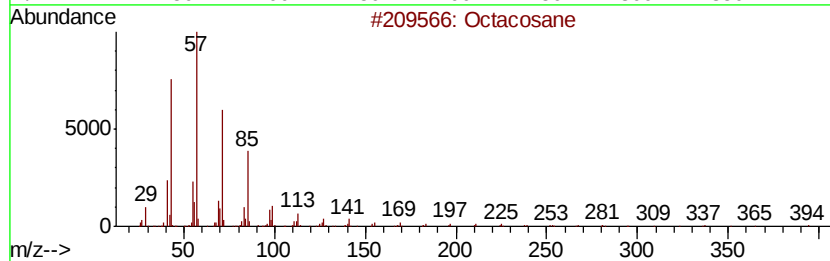
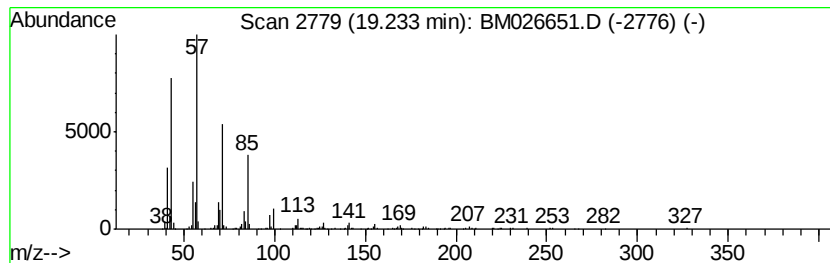
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.23	3.46 ng/ul	201494	Chrysene-d12	20.99

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octacosane	394	C28H58	000630-02-4	91
2		Heptacosane	380	C27H56	000593-49-7	90
3		Eicosane	282	C20H42	000112-95-8	87
4		Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	86
5		Tridecane, 7-hexyl-	268	C19H40	007225-66-3	86



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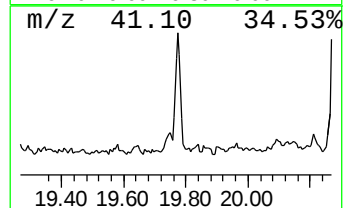
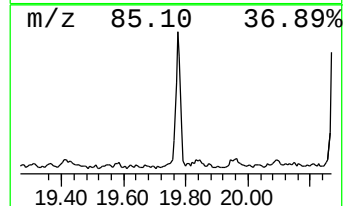
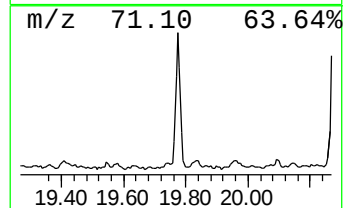
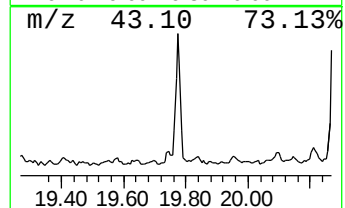
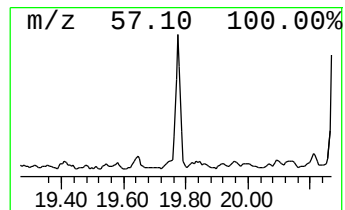
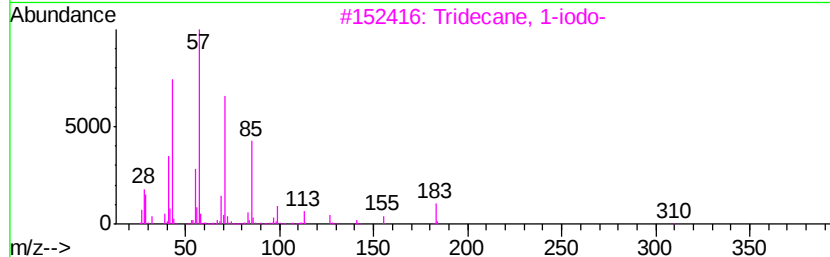
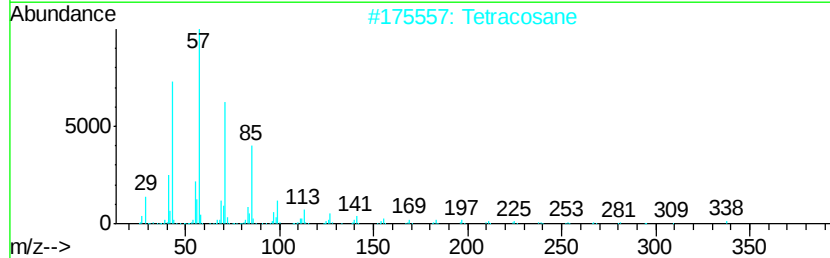
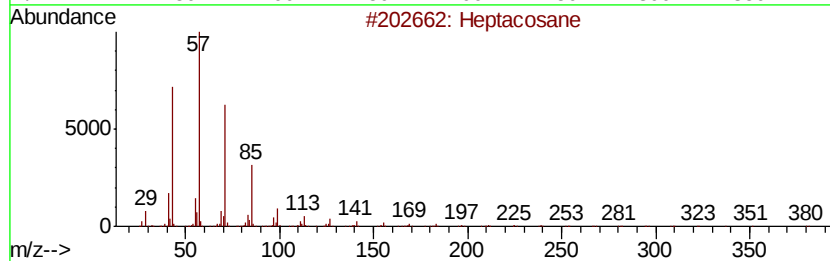
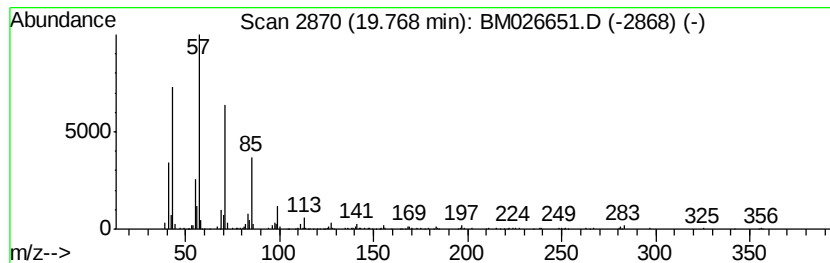
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.77	5.27 ng/ul	307089	Chrysene-d12	20.99

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptacosane	380	C27H56	000593-49-7	91
2		Tetracosane	338	C24H50	000646-31-1	91
3		Tridecane, 1-iodo-	310	C13H27I	035599-77-0	87
4		Docosane	310	C22H46	000629-97-0	86
5		Octadecane	254	C18H38	000593-45-3	86



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ALS Vial : 6 Sample Multiplier: 1

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ClientSampleId :
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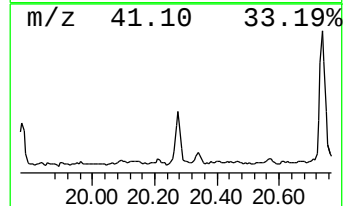
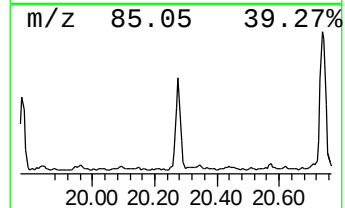
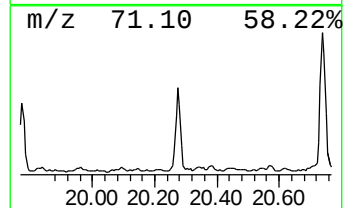
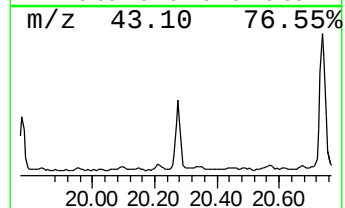
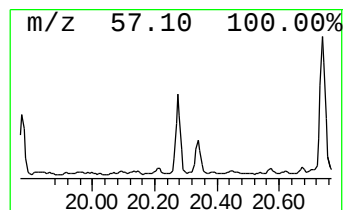
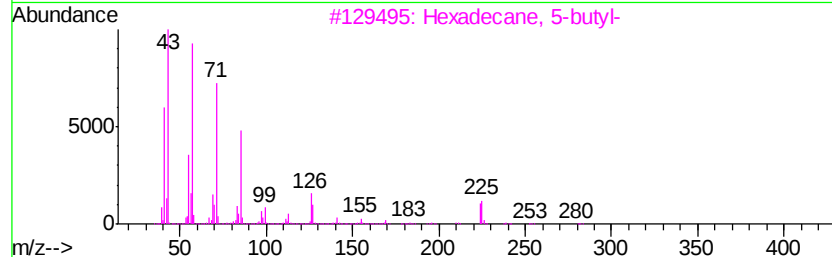
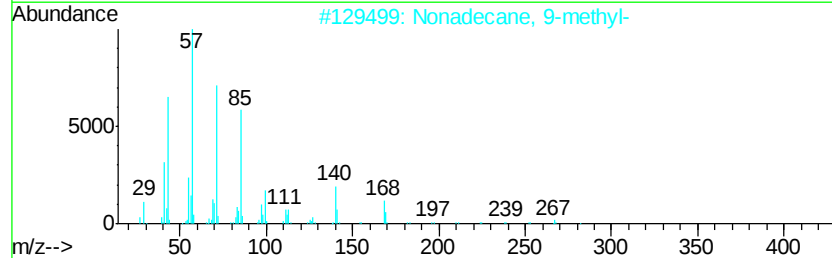
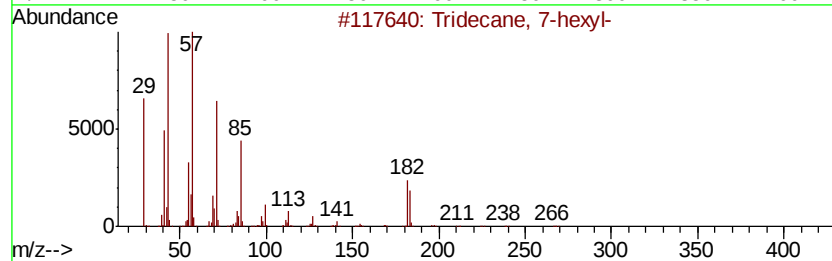
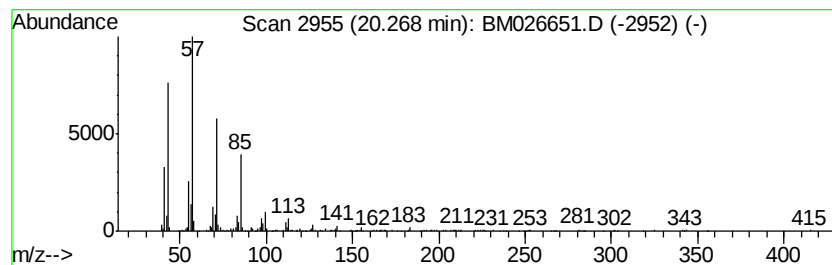
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.27	7.46 ng/ul	434928	Chrysene-d12	20.99

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tridecane, 7-hexyl-	268	C19H40	007225-66-3	95
2		Nonadecane, 9-methyl-	282	C20H42	013287-24-6	93
3		Hexadecane, 5-butyl-	282	C20H42	006912-07-8	93
4		Heneicosane	296	C21H44	000629-94-7	91
5		Heptacosane	380	C27H56	000593-49-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM070220\
Data File : BM026651.D
Acq On : 02 Jul 2020 13:08
Operator : JU/CG
Sample : L3131-01
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BG2L6

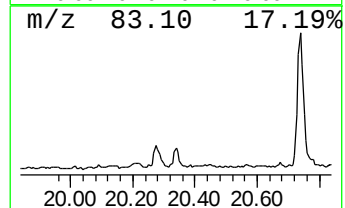
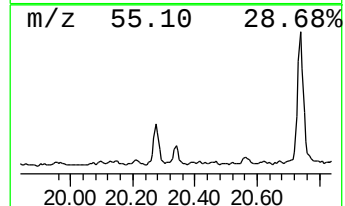
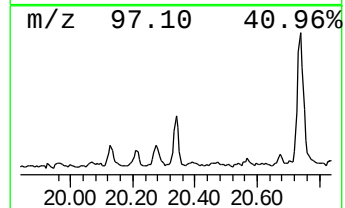
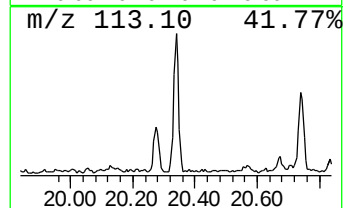
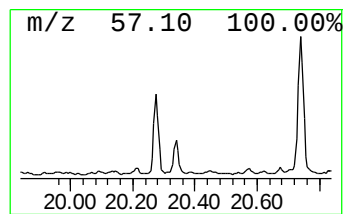
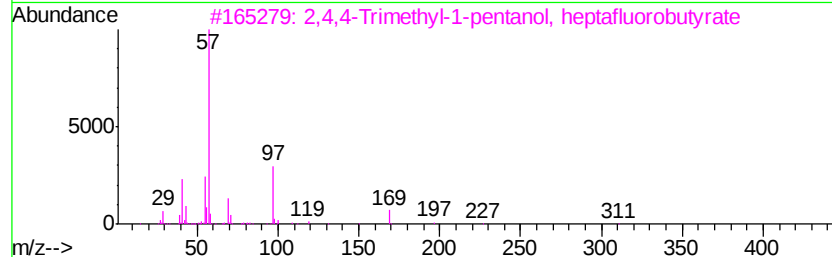
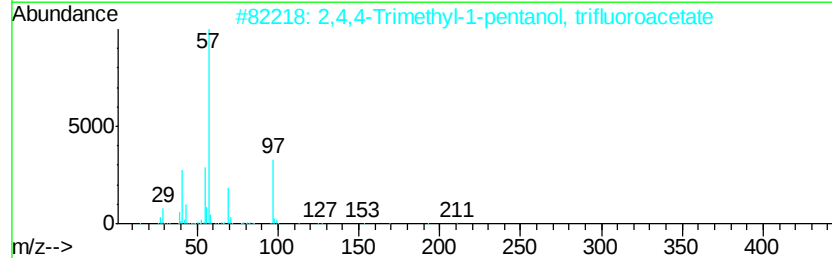
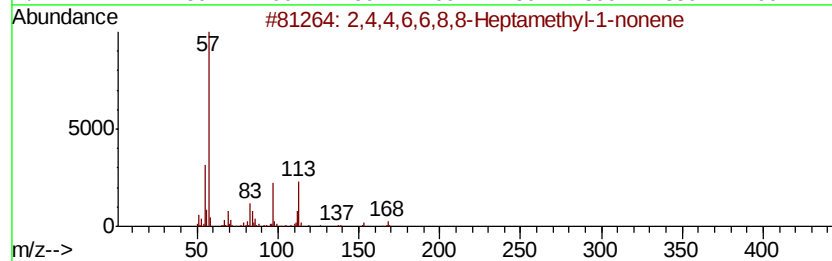
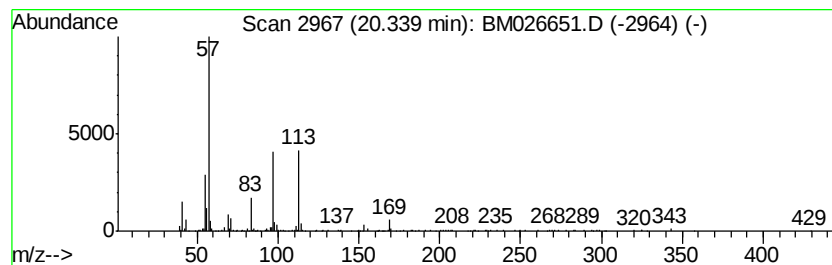
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.34	2.33 ng/ul	135938	Chrysene-d12	20.99

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	64		
2	2,4,4-Trimethyl-1-pentanol, trif...	226	C10H17F3O2	1000365-19-5	38		
3	2,4,4-Trimethyl-1-pentanol, hept...	326	C12H17F7O2	1000365-01-4	32		
4	2,4,4-Trimethyl-1-pentanol, pent...	276	C11H17F5O2	1000365-51-0	25		
5	1-Pentanol, 3-methyl-2-propyl-	144	C9H20O	054004-40-9	23		



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM070220\
Data File : BM026651.D
Acq On : 02 Jul 2020 13:08
Operator : JU/CG
Sample : L3131-01
Misc :
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Instrument :
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ClientSampleId :
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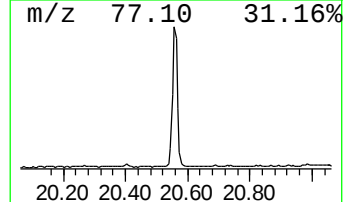
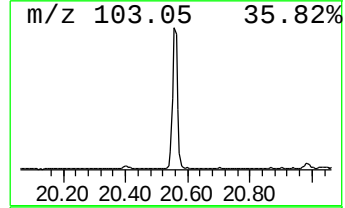
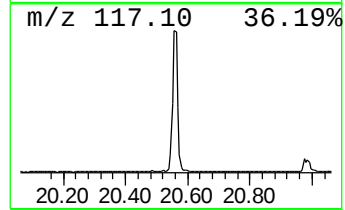
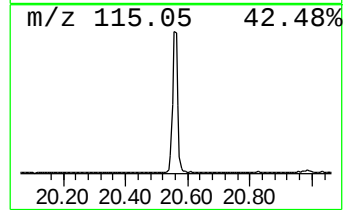
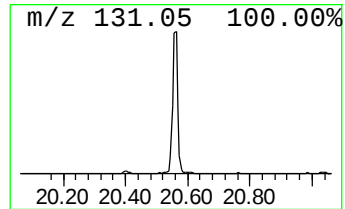
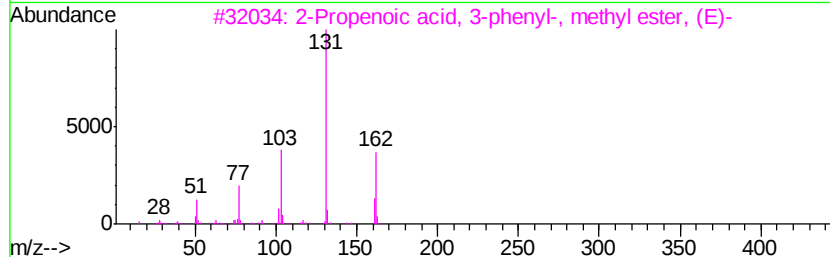
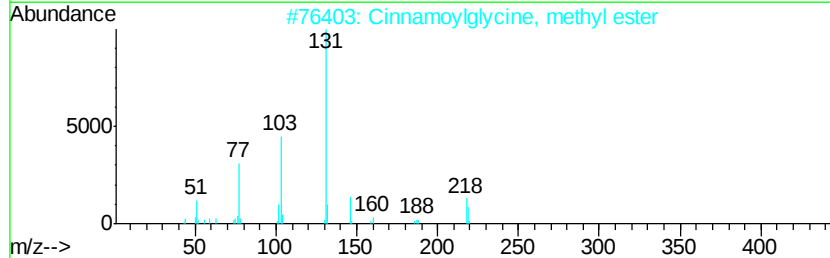
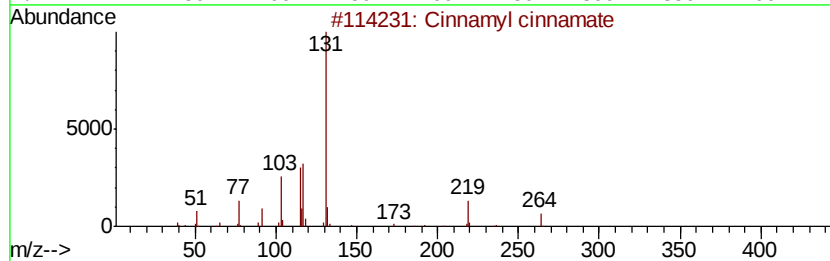
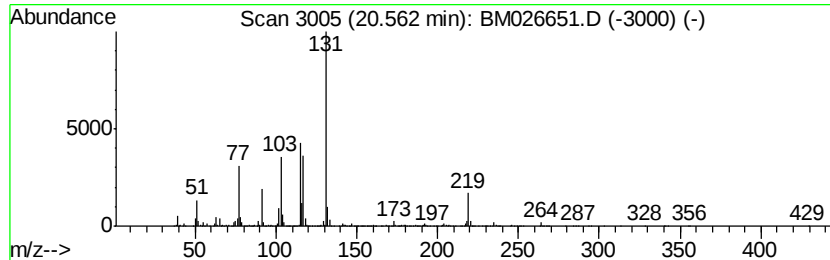
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 Cinnamyl cinnamate Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.56	16.50 ng/ul	961883	Chrysene-d12	20.99

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cinnamyl cinnamate	264	C18H16O2	000122-69-0	86
2		Cinnamoylglycine, methyl ester	219	C12H13NO3	040778-04-9	64
3		2-Propenoic acid, 3-phenyl-, met...	162	C10H10O2	001754-62-7	50
4		N-(p-Vinylbenzoyl)-l-alanine	219	C12H13NO3	139184-60-4	47
5		1-Penten-3-one, 4,4-dimethyl-1-p...	188	C13H16O	000538-44-3	47



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM070220\
Data File : BM026651.D
Acq On : 02 Jul 2020 13:08
Operator : JU/CG
Sample : L3131-01
Misc :
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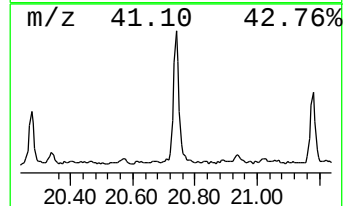
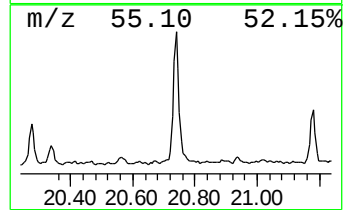
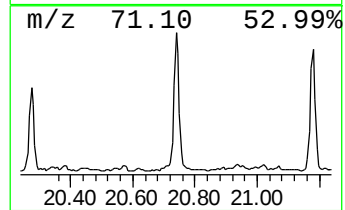
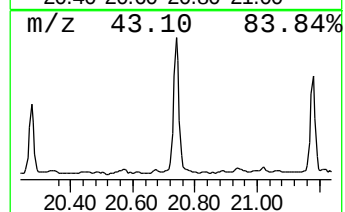
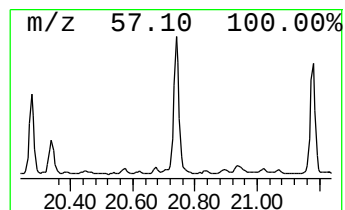
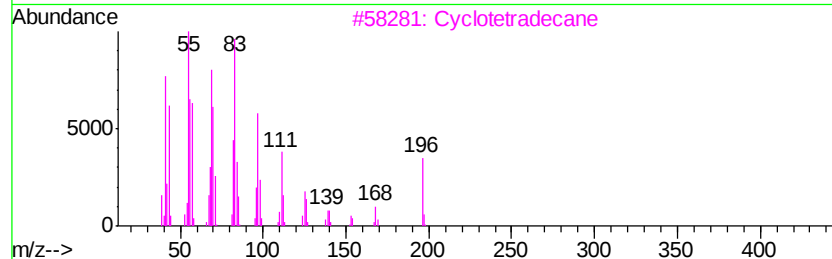
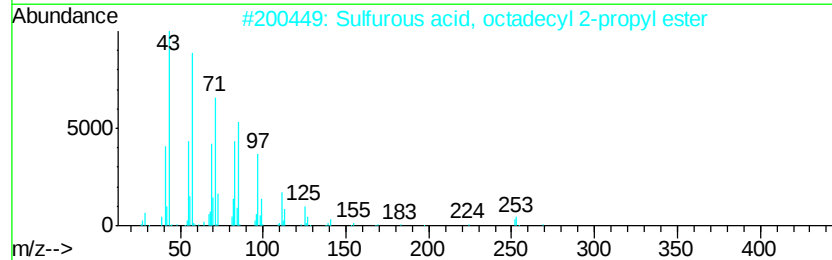
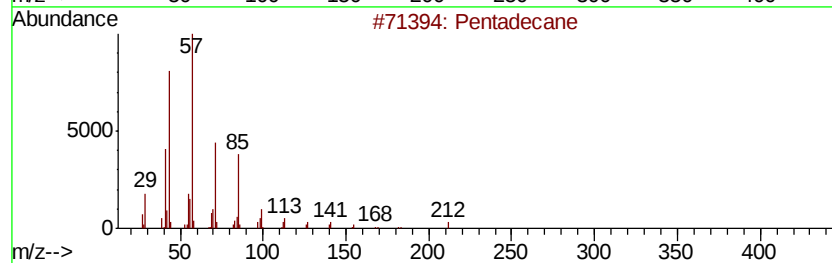
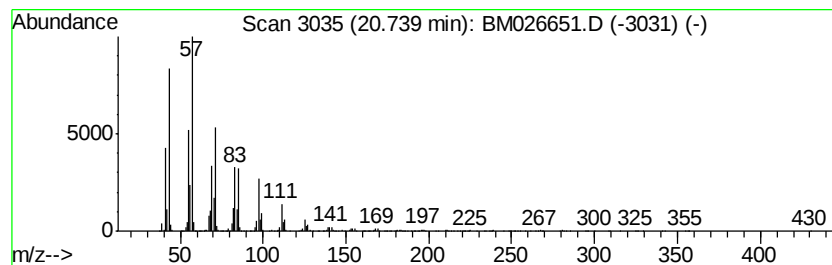
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.74	20.25 ng/ul	1180640	Chrysene-d12	20.99
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Pentadecane	212 C15H32	000629-62-9 90
2		Sulfurous acid, octadecyl 2-prop...	376 C21H44O3S	1000309-12-7 87
3		Cyclotetradecane	196 C14H28	000295-17-0 86
4		Oxalic acid, pentadecyl propyl e...	342 C20H38O4	1000309-26-8 83
5		Cyclohexadecane, 1,2-diethyl-	280 C20H40	1000155-85-3 78



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM070220\
Data File : BM026651.D
Acq On : 02 Jul 2020 13:08
Operator : JU/CG
Sample : L3131-01
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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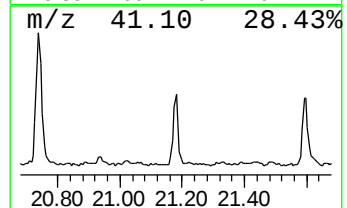
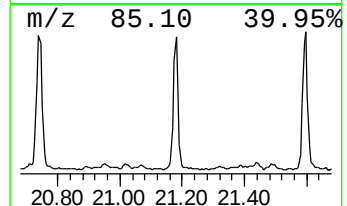
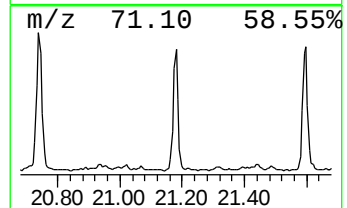
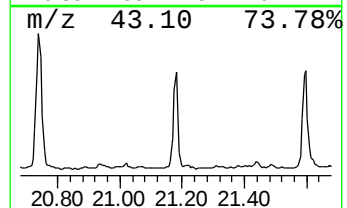
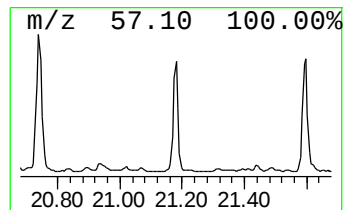
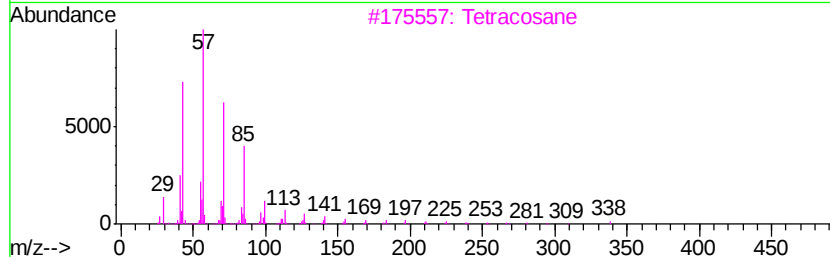
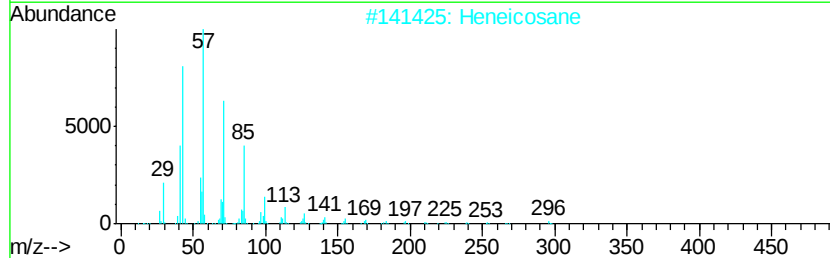
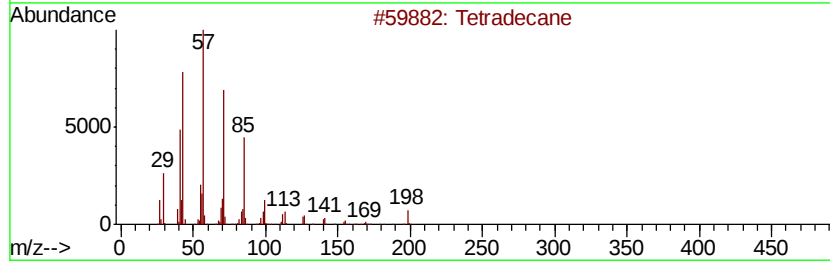
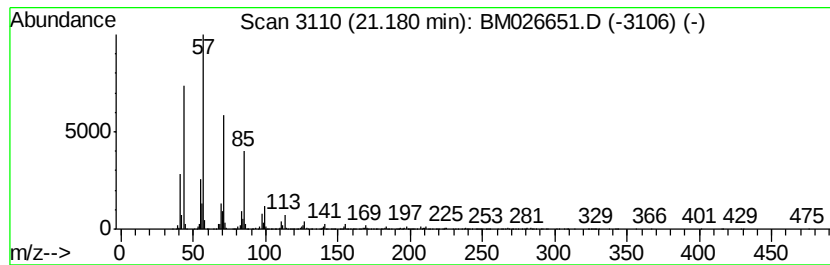
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.18	9.82 ng/ul	572688	Chrysene-d12	20.99

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetradecane	198	C14H30	000629-59-4	93
2			Heneicosane	296	C21H44	000629-94-7	91
3			Tetracosane	338	C24H50	000646-31-1	91
4			Heptacosane	380	C27H56	000593-49-7	91
5			Nonadecane	268	C19H40	000629-92-5	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM070220\
Data File : BM026651.D
Acq On : 02 Jul 2020 13:08
Operator : JU/CG
Sample : L3131-01
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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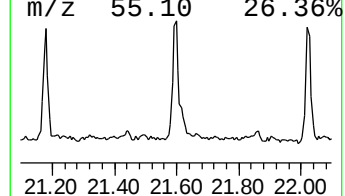
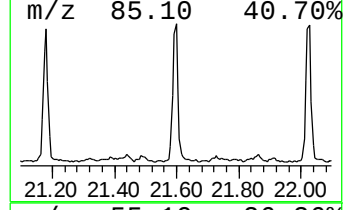
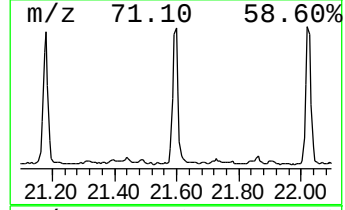
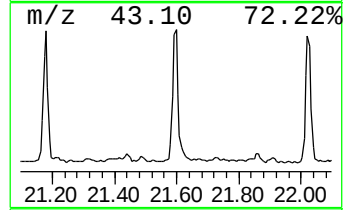
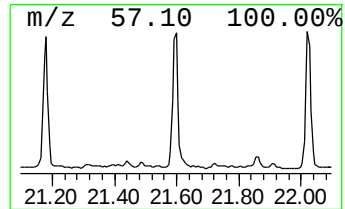
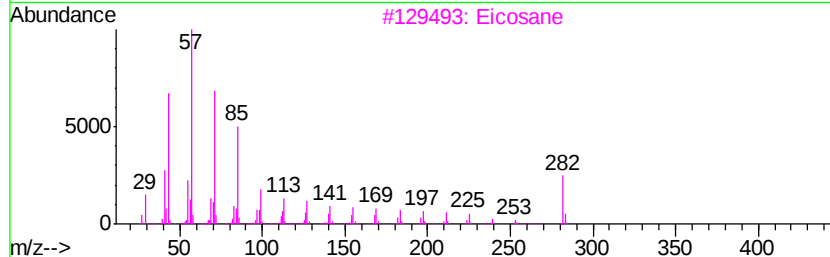
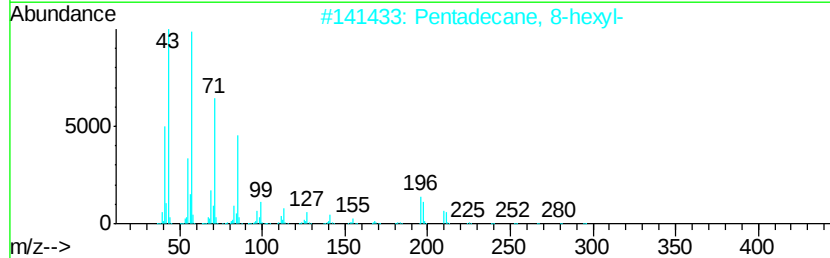
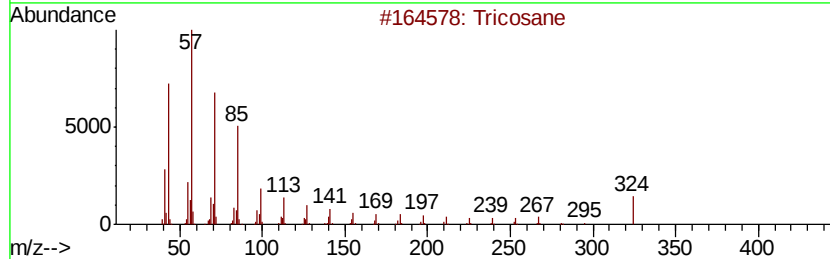
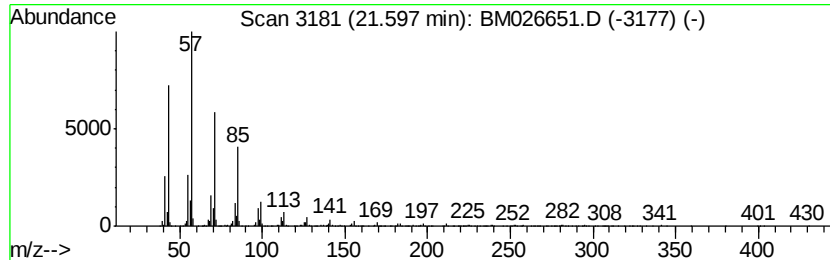
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Quant Title : SVOA CALIBRATION

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

Peak Number 15 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.60	12.29 ng/ul	716678	Chrysene-d12	20.99

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tricosane	324	C23H48	000638-67-5	97
2		Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	94
3		Eicosane	282	C20H42	000112-95-8	93
4		Tridecane, 7-hexyl-	268	C19H40	007225-66-3	93
5		Eicosane, 10-methyl-	296	C21H44	054833-23-7	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM070220\
Data File : BM026651.D
Acq On : 02 Jul 2020 13:08
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Sample : L3131-01
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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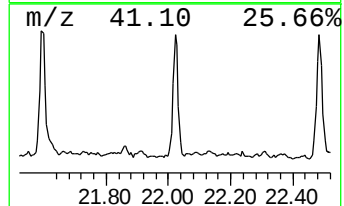
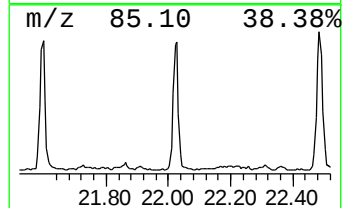
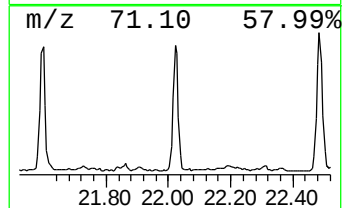
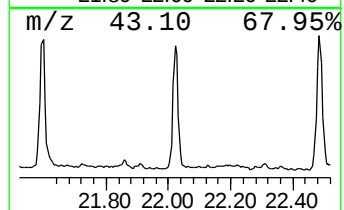
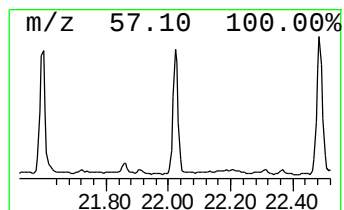
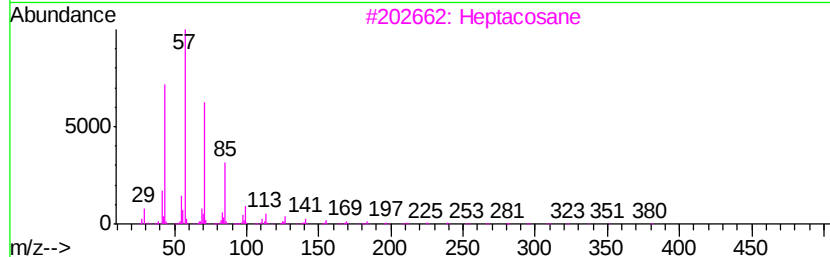
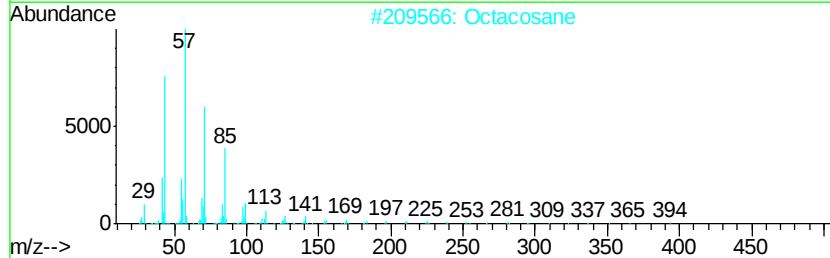
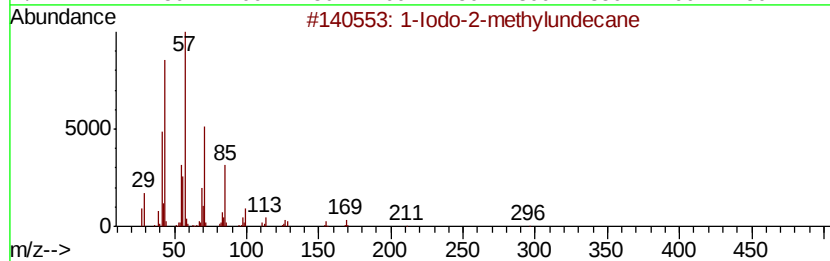
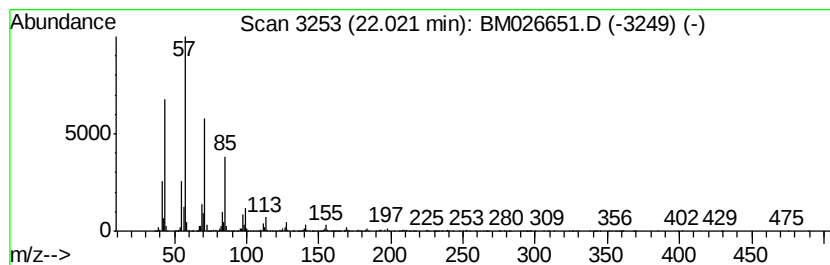
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 1-Iodo-2-methylundecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.02	11.39 ng/ul	664231	Chrysene-d12	20.99

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Iodo-2-methylundecane	296	C12H25I	073105-67-6	94
2		Octacosane	394	C28H58	000630-02-4	91
3		Heptacosane	380	C27H56	000593-49-7	91
4		Eicosane, 7-hexyl-	366	C26H54	055333-99-8	91
5		Heneicosane	296	C21H44	000629-94-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM070220\
Data File : BM026651.D
Acq On : 02 Jul 2020 13:08
Operator : JU/CG
Sample : L3131-01
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BG2L6

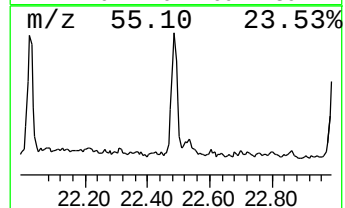
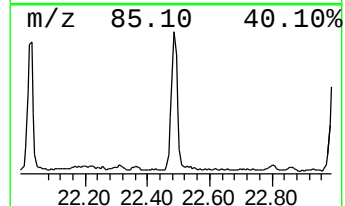
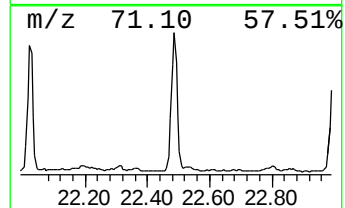
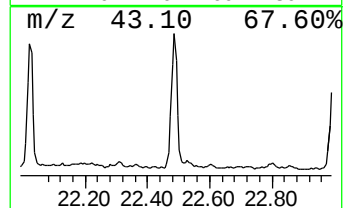
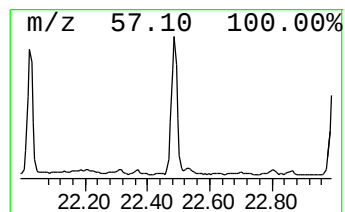
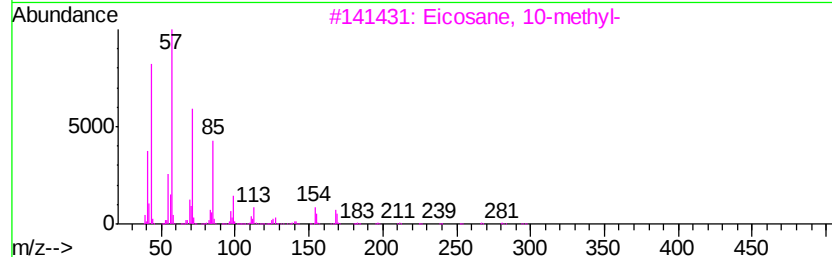
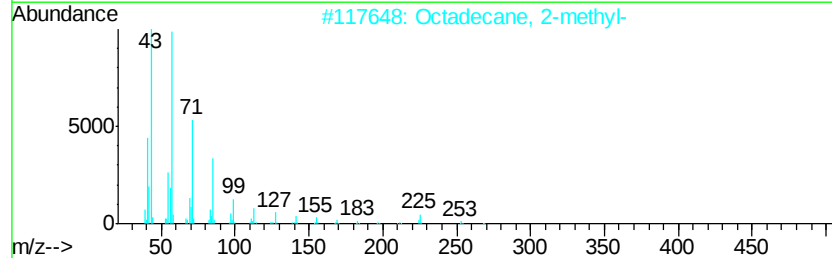
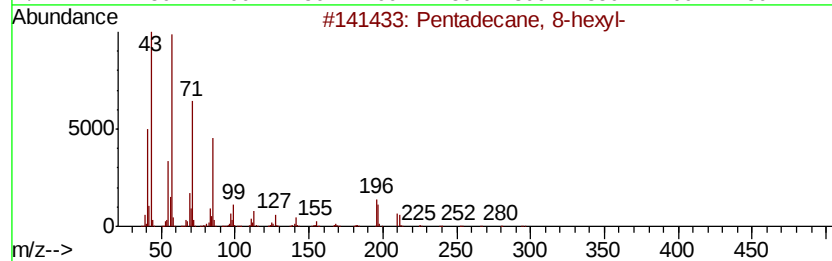
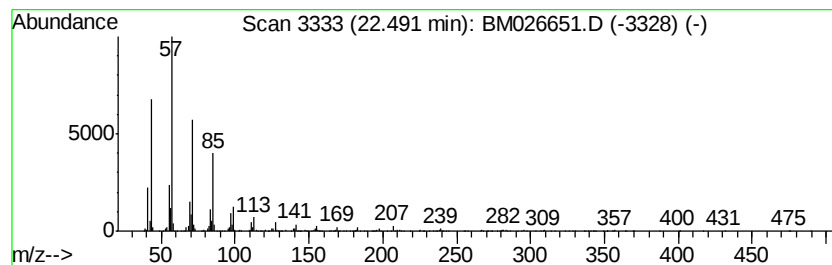
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Quant Title : SVOA CALIBRATION

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

Peak Number 17 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.49	11.80 ng/ul	730581	Perylene-d12	23.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	94
2		Octadecane, 2-methyl-	268	C19H40	001560-88-9	93
3		Eicosane, 10-methyl-	296	C21H44	054833-23-7	93
4		Octacosane	394	C28H58	000630-02-4	91
5		Tetratetracontane	619	C44H90	007098-22-8	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM070220\
Data File : BM026651.D
Acq On : 02 Jul 2020 13:08
Operator : JU/CG
Sample : L3131-01
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BG2L6

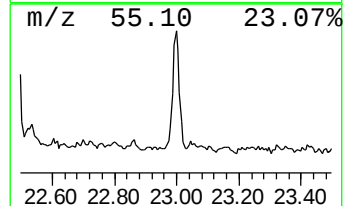
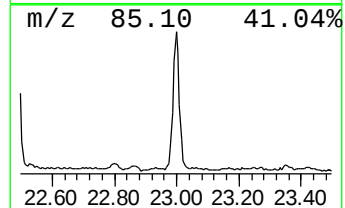
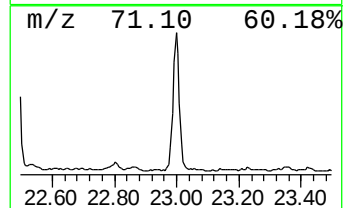
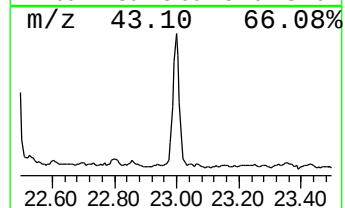
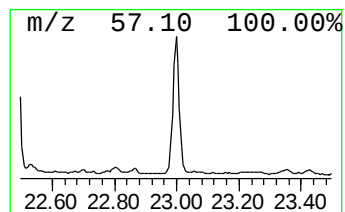
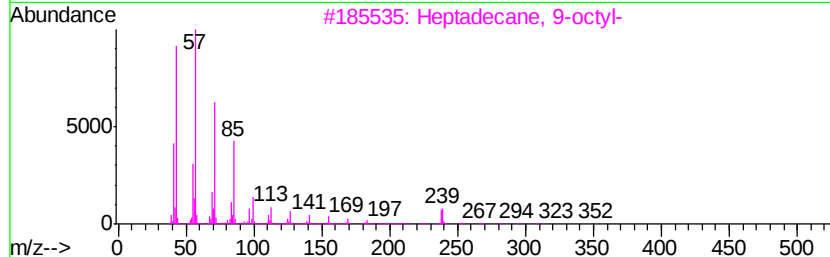
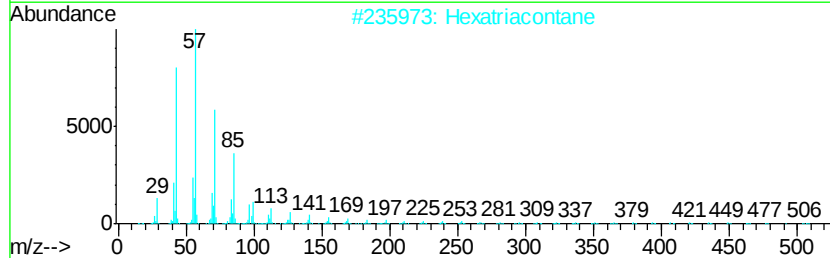
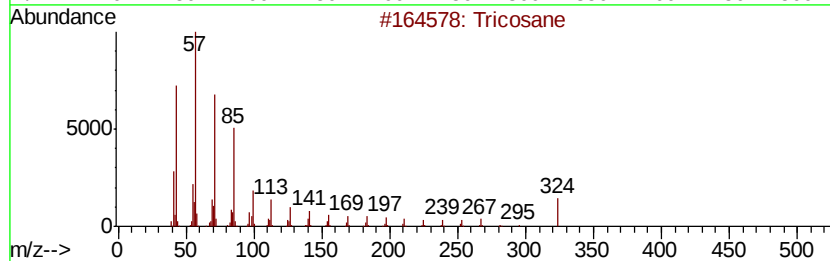
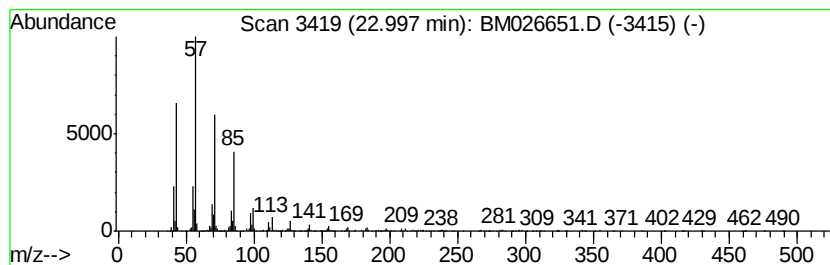
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Quant Title : SVOA CALIBRATION

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

Peak Number 18 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.00	10.76 ng/ul	666332	Perylene-d12	23.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tricosane	324	C23H48	000638-67-5	95
2		Hexatriacontane	507	C36H74	000630-06-8	91
3		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
4		Octacosane	394	C28H58	000630-02-4	91
5		Heptacosane	380	C27H56	000593-49-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM070220\
Data File : BM026651.D
Acq On : 02 Jul 2020 13:08
Operator : JU/CG
Sample : L3131-01
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BG2L6

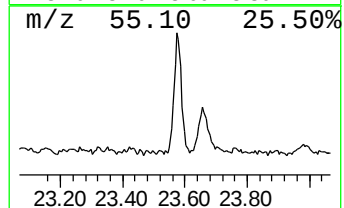
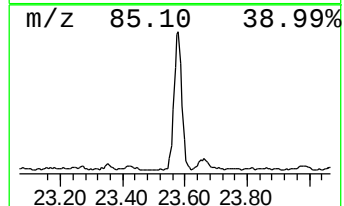
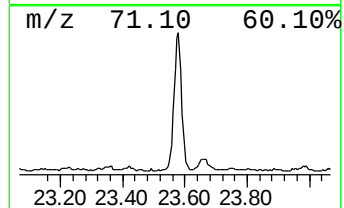
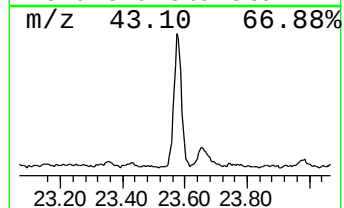
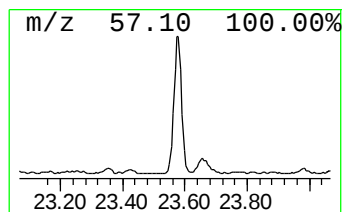
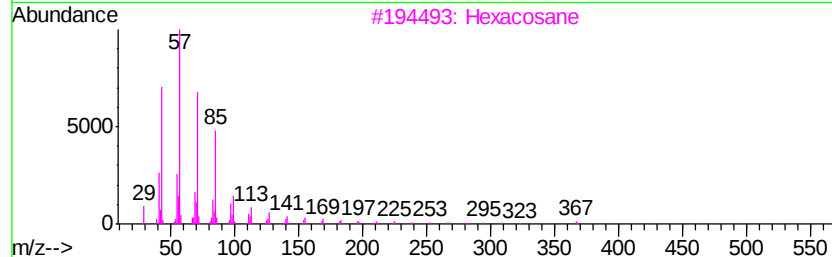
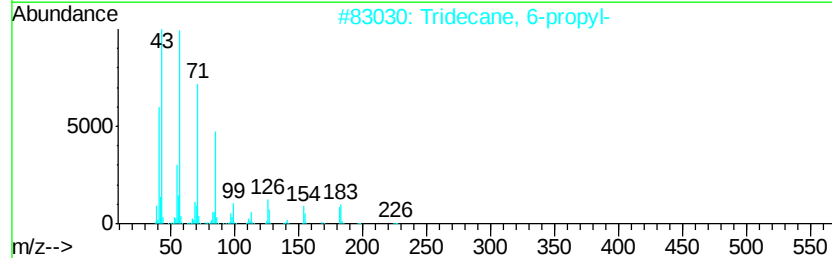
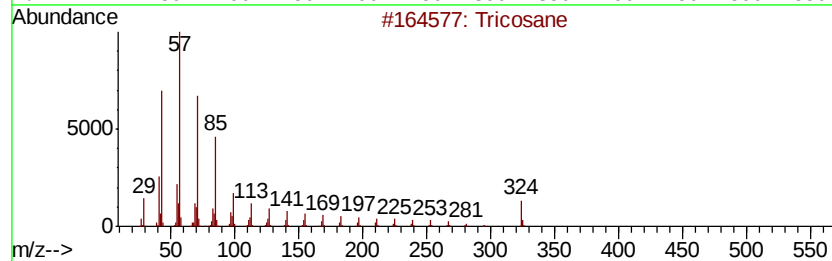
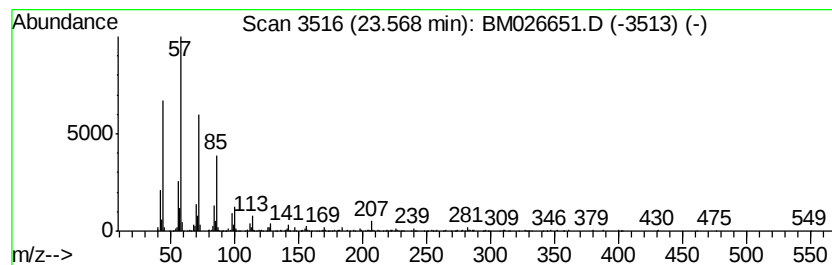
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Quant Title : SVOA CALIBRATION

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

Peak Number 19 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.57	11.40 ng/ul	705649	Perylene-d12	23.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tricosane	324	C23H48	000638-67-5	93
2		Tridecane, 6-propyl-	226	C16H34	055045-10-8	93
3		Hexacosane	366	C26H54	000630-01-3	91
4		Hexatriacontane	507	C36H74	000630-06-8	91
5		Octacosane	394	C28H58	000630-02-4	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM070220\
Data File : BM026651.D
Acq On : 02 Jul 2020 13:08
Operator : JU/CG
Sample : L3131-01
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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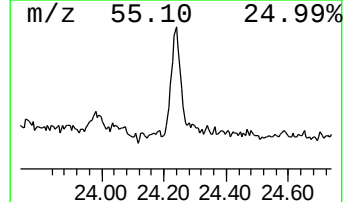
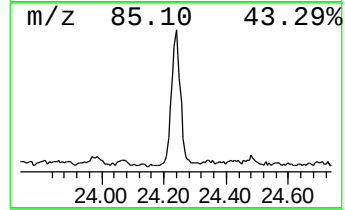
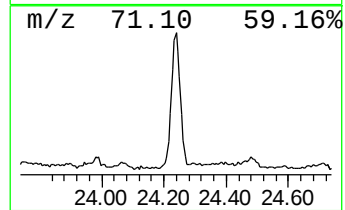
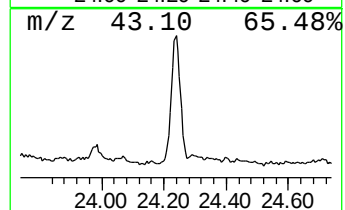
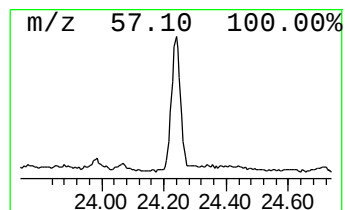
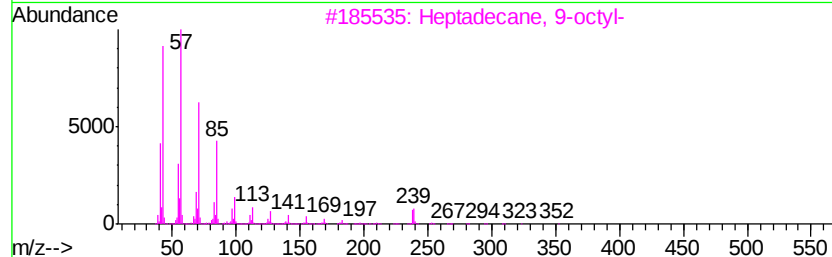
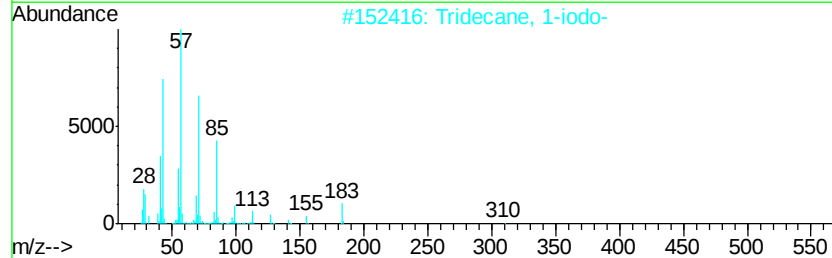
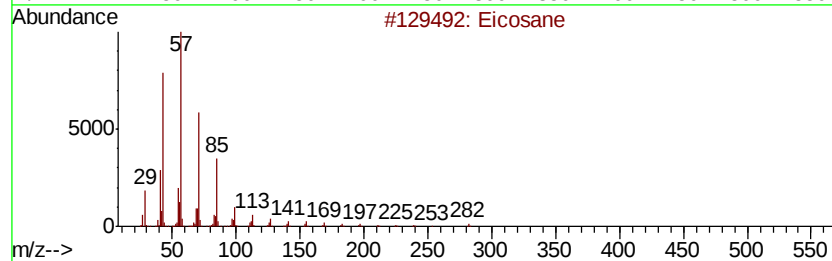
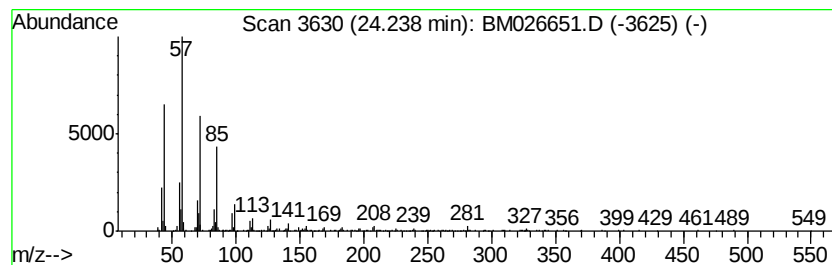
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Quant Title : SVOA CALIBRATION

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

Peak Number 20 (DEL) Alkane: Straight-Chai... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.24	6.19 ng/ul	383487	Perylene-d12	23.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	91
2		Tridecane, 1-iodo-	310	C13H27I	035599-77-0	91
3		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	90
4		Heptacosane	380	C27H56	000593-49-7	90
5		Hexacosane	366	C26H54	000630-01-3	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM070220\
Data File : BM026651.D
Acq On : 02 Jul 2020 13:08
Operator : JU/CG
Sample : L3131-01
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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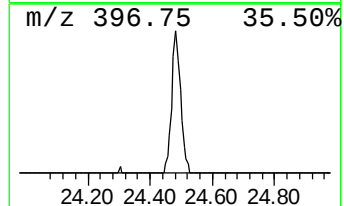
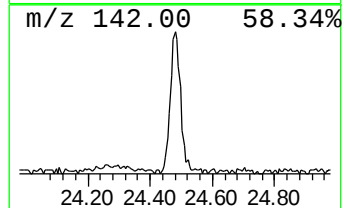
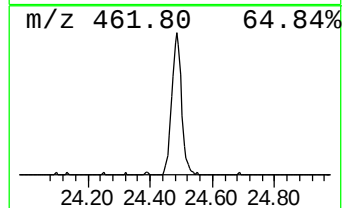
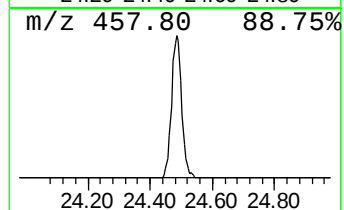
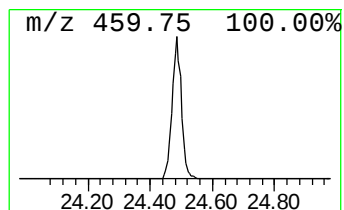
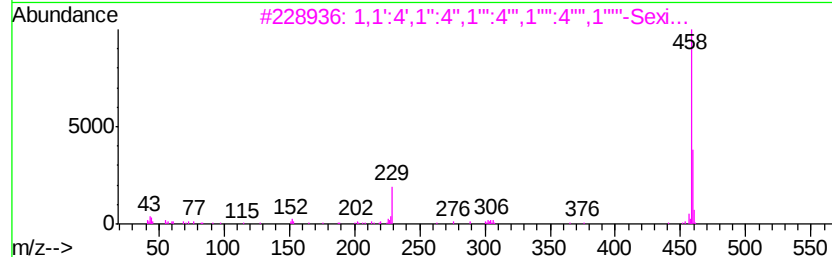
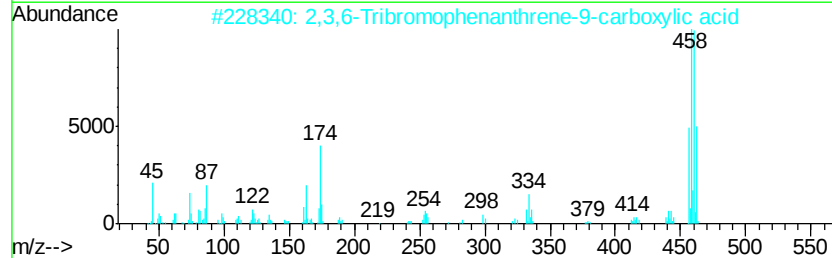
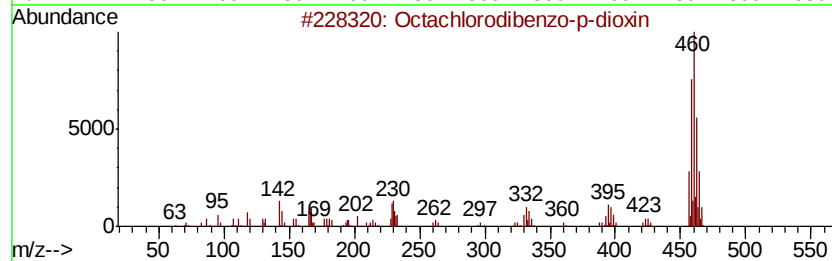
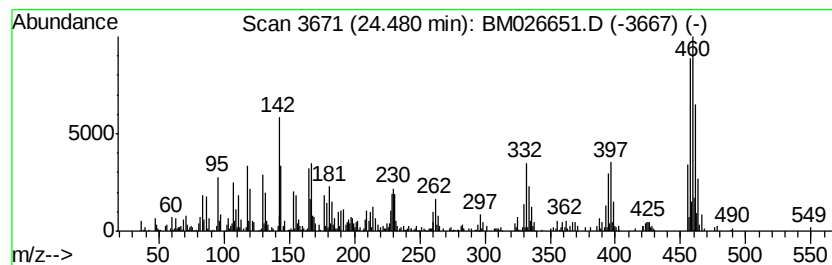
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Quant Title : SVOA CALIBRATION

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

Peak Number 21 Octachlorodibenzo-p-dioxin Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.48	8.61 ng/ul	533052	Perylene-d12	23.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octachlorodibenzo-p-dioxin	456	C12Cl8O2	003268-87-9	99
2		2,3,6-Tribromophenanthrene-9-car...	456	C15H7Br3O2	056988-60-4	46
3		1,1':4',1'':4'',1''':4''',1''':4''':...	458	C36H26	004499-83-6	14
4		(1-Amino-5-thiophen-2-yl-6,7,8,9...	458	C22H16Cl2N2O2S	1000318-00-6	10
5		Benzene, 1,4-bis(3-phenylindol-2...	460	C34H24N2	164936-88-3	10



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM070220\
 Data File : BM026651.D
 Acq On : 02 Jul 2020 13:08
 Operator : JU/CG
 Sample : L3131-01
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BG2L6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM063020MA.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
(1R)-2,6,6-Trimet...	6.06	3.2	ng/ul	101940	1	7.35	644716	20.0
Benzene, isocyanato-	6.43	2.4	ng/ul	77673	1	7.35	644716	20.0
2,4,7,9-Tetrameth...	12.93	6.5	ng/ul	311392	3	14.01	963290	20.0
n-Hexadecanoic acid	17.71	5.6	ng/ul	309720	4	16.77	1110750	20.0
(DEL) Alkane: Str...	18.46	4.3	ng/ul	236742	4	16.77	1110750	20.0
(DEL) Alkane: Str...	18.64	2.4	ng/ul	130319	4	16.77	1110750	20.0
Octadecanoic acid	19.01	3.0	ng/ul	171941	5	20.99	1166180	20.0
(DEL) Alkane: Str...	19.23	3.5	ng/ul	201494	5	20.99	1166180	20.0
(DEL) Alkane: Str...	19.77	5.3	ng/ul	307089	5	20.99	1166180	20.0
(DEL) Alkane: Str...	20.27	7.5	ng/ul	434928	5	20.99	1166180	20.0
(DEL) Alkane: Str...	20.34	2.3	ng/ul	135938	5	20.99	1166180	20.0
Cinnamyl cinnamate	20.56	16.5	ng/ul	961883	5	20.99	1166180	20.0
(DEL) Alkane: Str...	20.74	20.3	ng/ul	1180640	5	20.99	1166180	20.0
(DEL) Alkane: Str...	21.18	9.8	ng/ul	572688	5	20.99	1166180	20.0
(DEL) Alkane: Str...	21.60	12.3	ng/ul	716678	5	20.99	1166180	20.0
1-Iodo-2-methylun...	22.02	11.4	ng/ul	664231	5	20.99	1166180	20.0
(DEL) Alkane: Str...	22.49	11.8	ng/ul	730581	6	23.06	1238510	20.0
(DEL) Alkane: Str...	23.00	10.8	ng/ul	666332	6	23.06	1238510	20.0
(DEL) Alkane: Str...	23.57	11.4	ng/ul	705649	6	23.06	1238510	20.0
(DEL) Alkane: Str...	24.24	6.2	ng/ul	383487	6	23.06	1238510	20.0
Octachlorodibenzo...	24.48	8.6	ng/ul	533052	6	23.06	1238510	20.0