

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121721\
 Data File : BG051638.D
 Acq On : 18 Dec 2021 20:45
 Operator : CG/JU
 Sample : M5153-10
 Misc :
 ALS Vial : 54 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BGL75

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121421.M
 Title : SVOA CALIBRATION

Signal : TIC: BG051638.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.600	14	19	25	rBV3	7874	12386	0.59%	0.084%
2	4.029	88	92	108	rBV	52944	91279	4.36%	0.623%
3	7.154	617	624	630	rVB4	6018	13644	0.65%	0.093%
4	7.254	637	641	646	rVB2	7457	12100	0.58%	0.083%
5	7.401	660	666	676	rVB	143744	267223	12.77%	1.823%
6	7.577	689	696	707	rBV	100448	180066	8.61%	1.228%
7	7.800	723	734	746	rVB	129521	253082	12.10%	1.726%
8	8.188	797	800	806	rVV4	7942	16052	0.77%	0.109%
9	8.270	806	814	820	rVV2	128173	258809	12.37%	1.765%
10	8.346	822	827	833	rVB5	9899	21664	1.04%	0.148%
11	8.470	842	848	853	rBV2	30334	53136	2.54%	0.362%
12	8.529	853	858	862	rVV3	12359	26199	1.25%	0.179%
13	8.570	862	865	870	rVB3	9937	14269	0.68%	0.097%
14	8.963	926	932	940	rVV2	173772	312904	14.96%	2.134%
15	9.134	956	961	963	rBV5	9918	14352	0.69%	0.098%
16	9.439	1003	1013	1020	rVV	127060	253925	12.14%	1.732%
17	9.609	1037	1042	1045	rBV5	11083	19993	0.96%	0.136%
18	9.656	1045	1050	1056	rVB5	11163	23549	1.13%	0.161%
19	9.874	1081	1087	1091	rVB6	11409	23840	1.14%	0.163%
20	9.938	1091	1098	1101	rBV5	15221	26226	1.25%	0.179%
21	10.173	1130	1138	1144	rBV	115880	219449	10.49%	1.497%
22	10.226	1144	1147	1151	rVB3	9056	12818	0.61%	0.087%
23	10.291	1151	1158	1163	rBV8	7475	13111	0.63%	0.089%
24	10.714	1221	1230	1238	rVV	171148	324839	15.53%	2.216%
25	11.113	1286	1298	1304	rBV	164539	316486	15.13%	2.159%
26	11.231	1312	1318	1330	rVB4	103238	240614	11.50%	1.641%
27	11.942	1434	1439	1444	rBV6	7178	11990	0.57%	0.082%
28	12.024	1448	1453	1462	rVB6	5594	15011	0.72%	0.102%
29	12.124	1462	1470	1473	rBV3	32476	56549	2.70%	0.386%
30	12.500	1528	1534	1538	rBV6	8231	13579	0.65%	0.093%
31	12.729	1566	1573	1580	rVB4	31079	62067	2.97%	0.423%
32	12.858	1587	1595	1599	rVB8	7444	14567	0.70%	0.099%
33	12.970	1608	1614	1624	rVB10	7983	22729	1.09%	0.155%
34	13.134	1637	1642	1645	rBV5	9822	16165	0.77%	0.110%
35	13.445	1690	1695	1701	rVB2	44336	63936	3.06%	0.436%
36	13.722	1736	1742	1751	rBV7	17899	53115	2.54%	0.362%

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37	13.816	1751	1758	1760	rVV3	11916	25180	1.20%	0.172%
38	13.851	1762	1764	1769	rVV5	11437	20119	0.96%	0.137%
39	13.892	1769	1771	1776	rVV6	8196	11721	0.56%	0.080%
40	13.968	1781	1784	1789	rVB4	9067	11885	0.57%	0.081%
41	14.191	1817	1822	1825	rBV7	6186	13168	0.63%	0.090%
42	14.238	1826	1830	1834	rVV7	14329	24891	1.19%	0.170%
43	14.297	1834	1840	1845	rVV	307743	450959	21.56%	3.076%
44	14.344	1845	1848	1851	rVV2	22159	32508	1.55%	0.222%
45	14.379	1851	1854	1859	rVB2	47890	67347	3.22%	0.459%
46	14.491	1871	1873	1877	rVB5	9804	12041	0.58%	0.082%
47	14.556	1878	1884	1886	rBV7	8310	13407	0.64%	0.091%
48	14.603	1886	1892	1899	rBV	344565	545450	26.07%	3.721%
49	14.750	1910	1917	1920	rBV5	22906	36804	1.76%	0.251%
50	14.797	1921	1925	1933	rVV6	13967	37727	1.80%	0.257%
51	14.908	1933	1944	1951	rVB	274483	482786	23.08%	3.293%
52	14.990	1956	1958	1963	rVV5	13755	16078	0.77%	0.110%
53	15.067	1965	1971	1977	rVV	152806	246938	11.80%	1.684%
54	15.126	1979	1981	1988	rVB8	13626	20030	0.96%	0.137%
55	15.231	1995	1999	2001	rBV5	12644	17591	0.84%	0.120%
56	15.290	2005	2009	2016	rVB8	23646	44274	2.12%	0.302%
57	15.419	2026	2031	2036	rBV8	18281	40317	1.93%	0.275%
58	15.484	2036	2042	2044	rBV5	24043	44692	2.14%	0.305%
59	15.531	2046	2050	2055	rVB7	23228	43630	2.09%	0.298%
60	15.631	2062	2067	2070	rBV7	11907	20551	0.98%	0.140%
61	15.689	2074	2077	2090	rVB7	35731	101773	4.86%	0.694%
62	15.789	2092	2094	2099	rBV6	9443	16960	0.81%	0.116%
63	15.842	2100	2103	2107	rBV5	16376	22888	1.09%	0.156%
64	15.895	2107	2112	2118	rVV	446270	676483	32.34%	4.614%
65	15.995	2124	2129	2134	rBV2	118324	176856	8.45%	1.206%
66	16.083	2141	2144	2147	rVB2	17817	20559	0.98%	0.140%
67	16.148	2151	2155	2159	rVB	45199	68536	3.28%	0.467%
68	16.236	2167	2170	2177	rVB9	15377	30736	1.47%	0.210%
69	16.359	2188	2191	2194	rVB4	15652	17155	0.82%	0.117%
70	16.394	2194	2197	2201	rBV2	23937	25564	1.22%	0.174%
71	16.623	2228	2236	2241	rBV2	121760	210388	10.06%	1.435%
72	16.723	2249	2253	2256	rBV	67471	90861	4.34%	0.620%
73	16.776	2258	2262	2268	rVV5	65343	125320	5.99%	0.855%
74	16.841	2269	2273	2276	rVV2	56881	77414	3.70%	0.528%
75	16.882	2276	2280	2283	rVB4	44147	65699	3.14%	0.448%
76	17.046	2303	2308	2314	rVB3	56900	120710	5.77%	0.823%

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77	17.111	2314	2319	2321	rBV	57547	74602	3.57%	0.509%
78	17.182	2328	2331	2335	rBV3	27281	34878	1.67%	0.238%
79	17.446	2373	2376	2381	rVB	50092	72814	3.48%	0.497%
80	17.540	2389	2392	2398	rVB3	41530	64632	3.09%	0.441%
81	17.657	2406	2412	2417	rBV	345743	526370	25.16%	3.590%
82	17.751	2423	2428	2435	rVB2	465184	715404	34.20%	4.880%
83	18.057	2477	2480	2486	rVB6	18687	30607	1.46%	0.209%
84	18.304	2518	2522	2527	rVB7	31651	48932	2.34%	0.334%
85	18.527	2556	2560	2564	rVB5	37222	50193	2.40%	0.342%
86	19.091	2653	2656	2660	rVB5	42728	53485	2.56%	0.365%
87	19.431	2711	2714	2719	rVB7	32884	51783	2.48%	0.353%
88	19.625	2744	2747	2752	rVB	67251	91836	4.39%	0.626%
89	19.901	2792	2794	2800	rVB2	37692	47699	2.28%	0.325%
90	20.031	2811	2816	2822	rVB	549650	823816	39.38%	5.619%
91	21.587	3078	3081	3090	rVB4	49542	88930	4.25%	0.607%
92	21.846	3119	3125	3134	rVB	1431472	2092033	100.00%	14.270%
93	21.975	3141	3147	3154	rVB2	325251	590250	28.21%	4.026%
94	22.839	3291	3294	3299	rVB5	17481	29508	1.41%	0.201%
95	23.050	3328	3330	3334	rVB4	19560	22859	1.09%	0.156%
96	23.168	3345	3350	3356	rVB	46097	89492	4.28%	0.610%
97	23.602	3420	3424	3433	rVB7	22650	47668	2.28%	0.325%
98	24.883	3638	3642	3648	rBV2	27592	61349	2.93%	0.418%
99	25.235	3691	3702	3712	rVB2	287557	905429	43.28%	6.176%
100	25.476	3733	3743	3754	rBV2	196168	566089	27.06%	3.861%

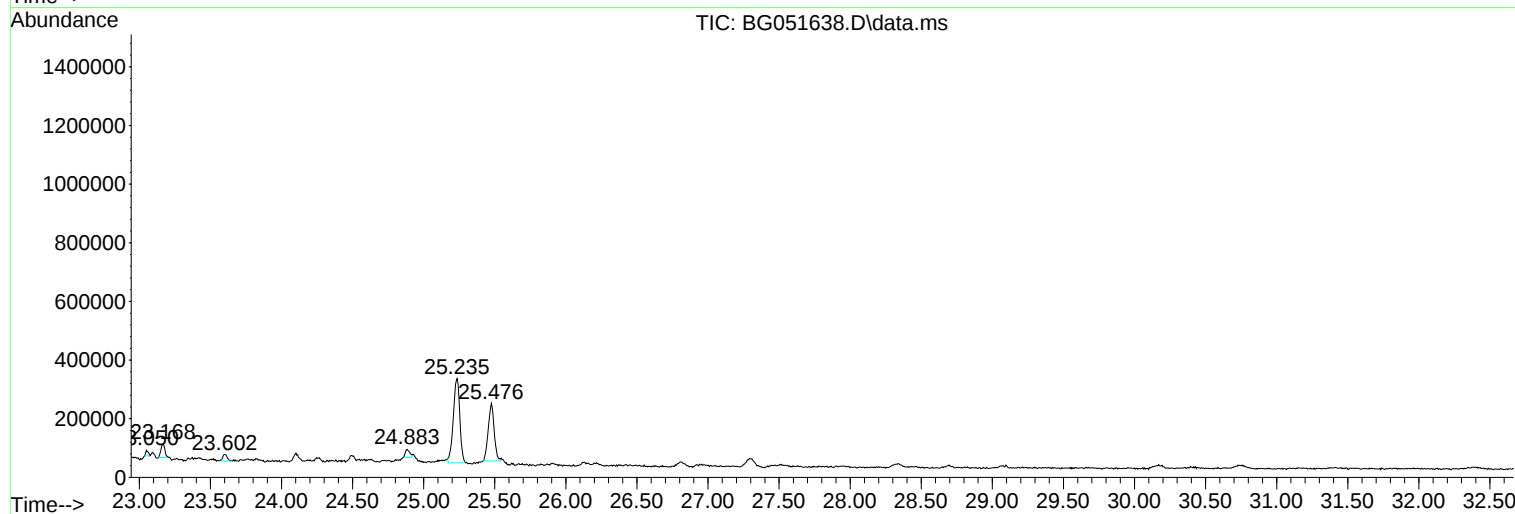
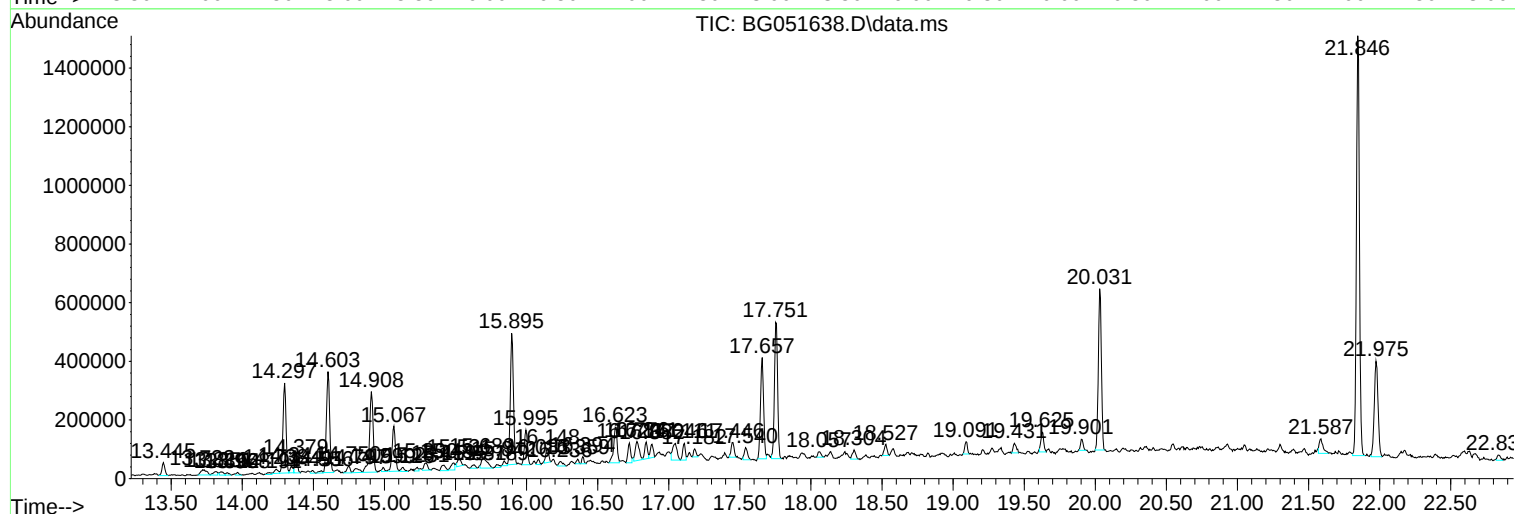
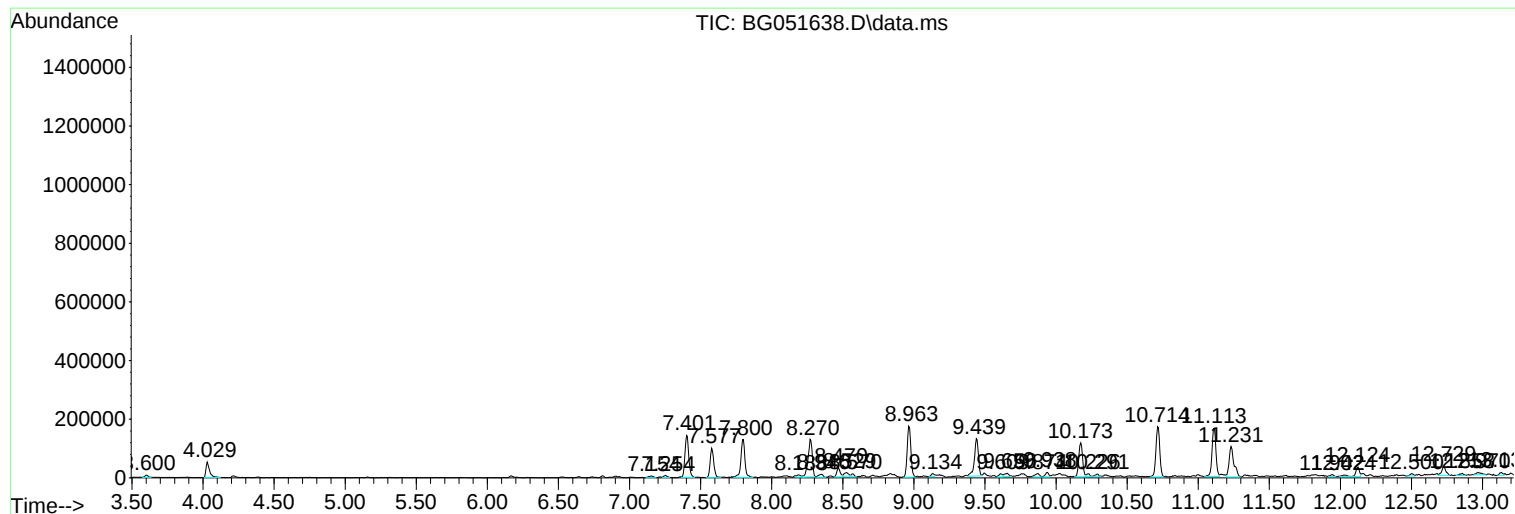
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121421.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P



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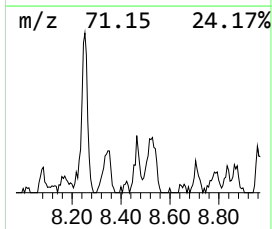
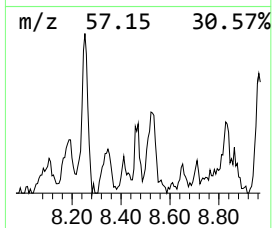
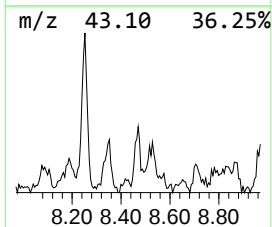
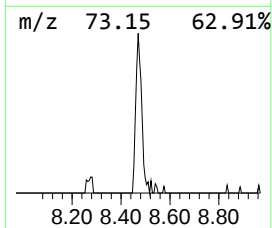
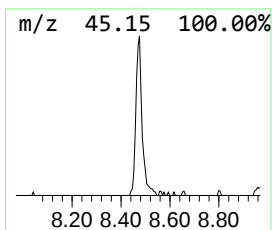
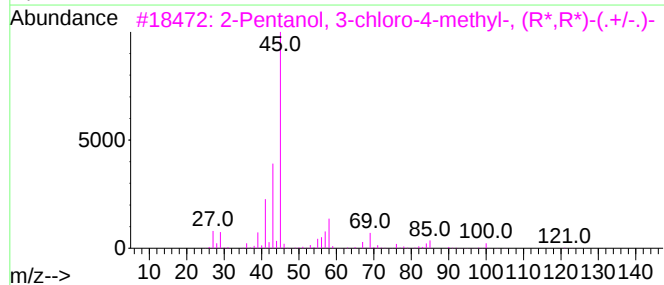
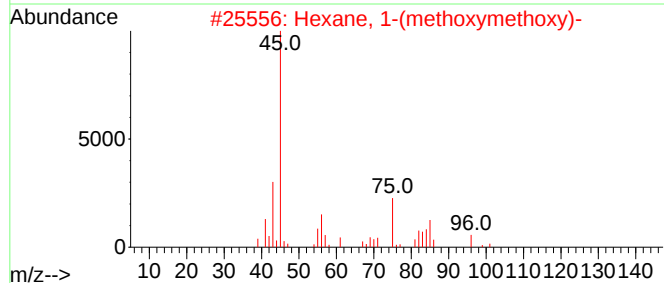
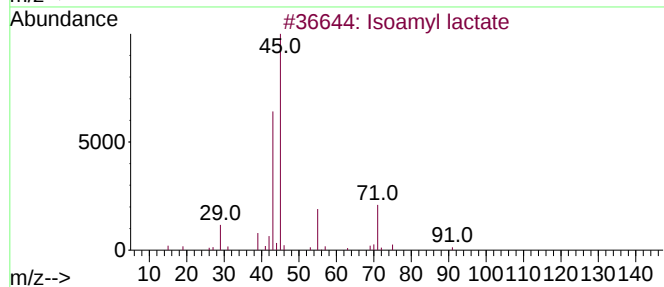
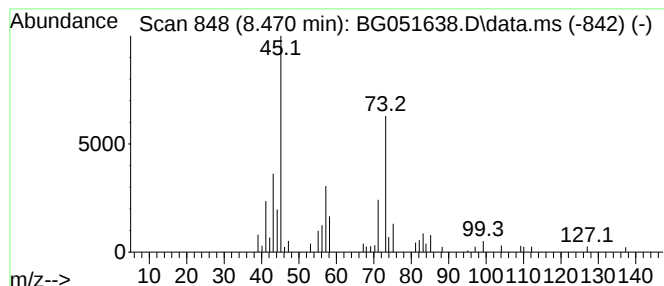
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121421.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 unknown-01 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.470	4.11 ng/ul	53136	1,4-Dichlorobenzene-d4	8.276

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Isoamyl lactate	160	C8H16O3	019329-89-6	47
2			Hexane, 1-(methoxymethoxy)-	146	C8H18O2	066675-06-7	47
3			2-Pentanol, 3-chloro-4-methyl-, ...	136	C6H13ClO	074685-47-5	43
4			Triethylene glycol	150	C6H14O4	000112-27-6	42
5			Ethanol, 2-[2-(2-butoxyethoxy)et...	206	C10H22O4	000143-22-6	40



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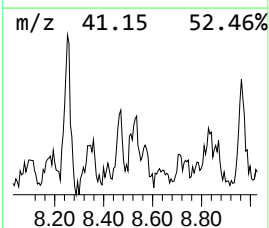
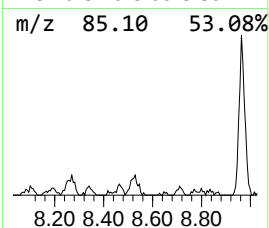
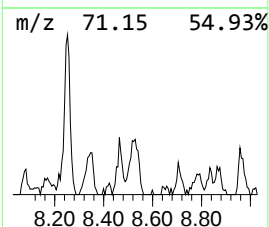
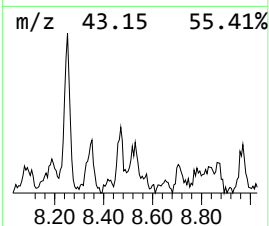
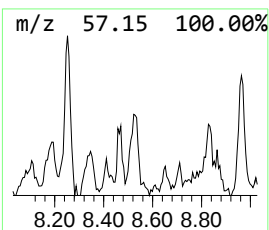
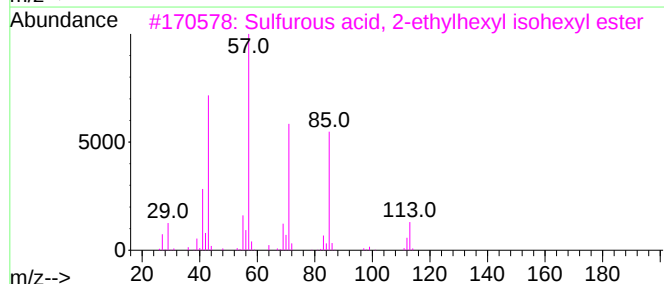
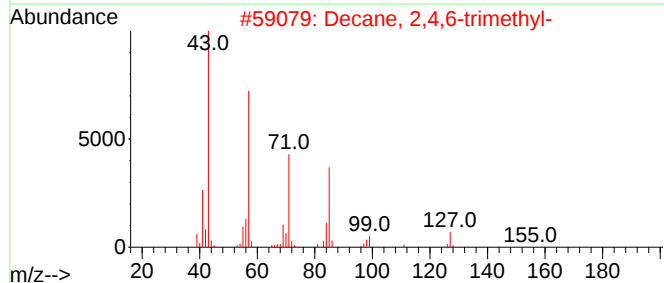
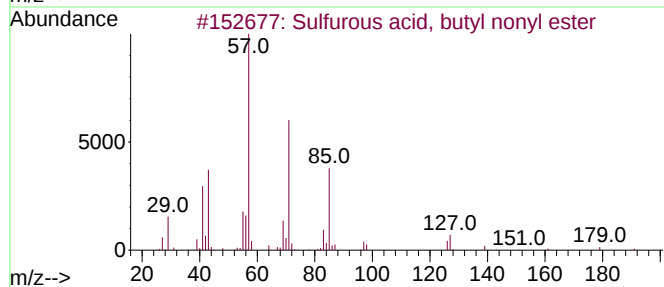
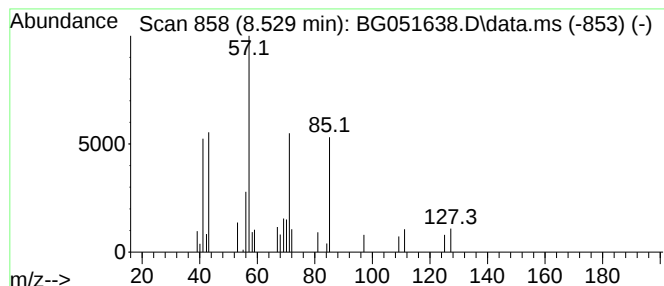
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Sulfurous acid, butyl nonyl... Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.529	2.02 ng/ul	26199	1,4-Dichlorobenzene-d4	8.276

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Sulfurous acid, butyl nonyl ester	264	C13H28O3S	1000309-17-6	53
2			Decane, 2,4,6-trimethyl-	184	C13H28	062108-27-4	50
3			Sulfurous acid, 2-ethylhexyl iso...	278	C14H30O3S	1000309-19-0	47
4			Carbonic acid, hexadecyl prop-1-...	326	C20H38O3	1000382-90-3	45
5			Carbonic acid, eicosyl vinyl ester	368	C23H44O3	1000382-54-3	45



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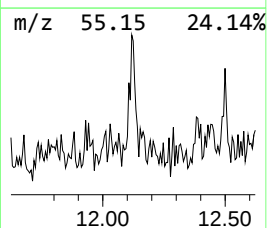
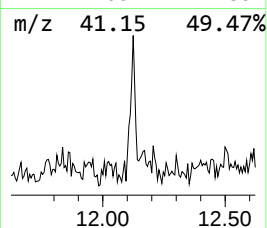
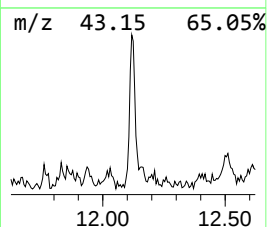
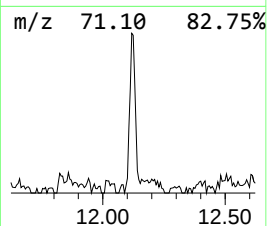
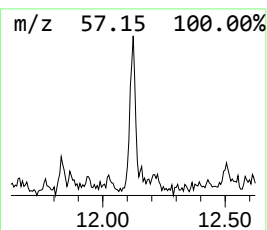
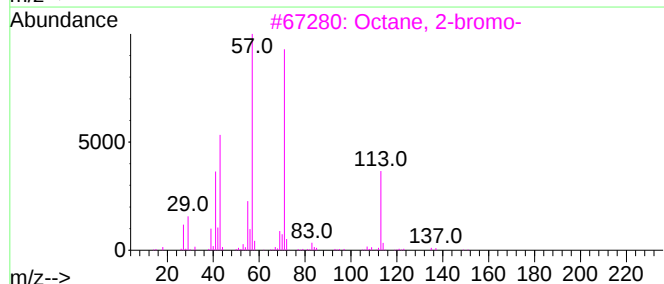
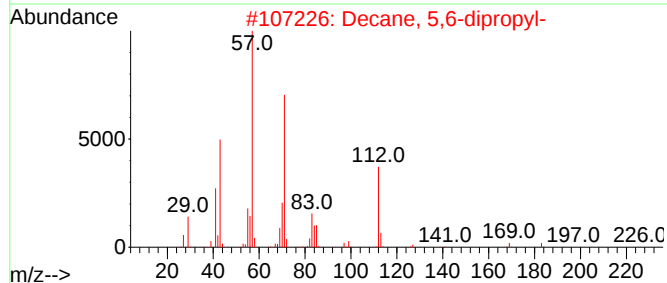
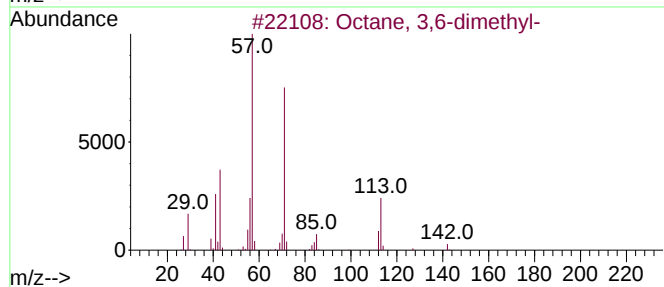
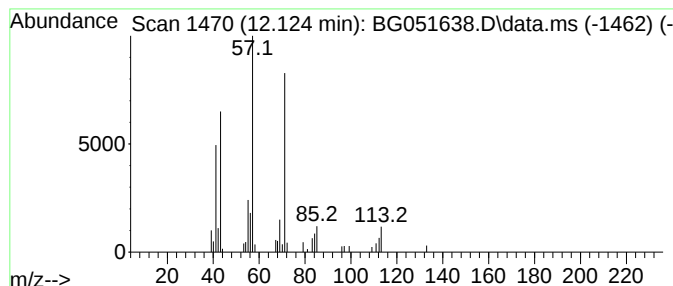
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121421.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.124	3.57 ng/ul	56549	Naphthalene-d8	11.113

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	78
2		Decane, 5,6-dipropyl-	226	C16H34	119209-20-0	78
3		Octane, 2-bromo-	192	C8H17Br	000557-35-7	72
4		Decane, 4-methyl-	156	C11H24	002847-72-5	64
5		3-Bromooctane	192	C8H17Br	000999-64-4	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121721\
Data File : BG051638.D
Acq On : 18 Dec 2021 20:45
Operator : CG/JU
Sample : M5153-10
Misc :
ALS Vial : 54 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGL75

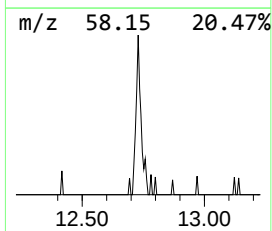
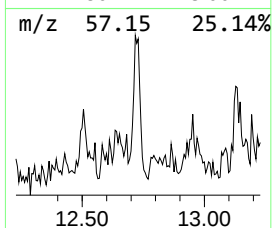
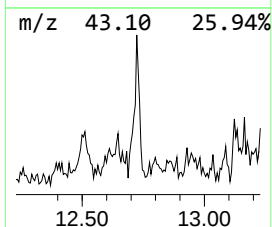
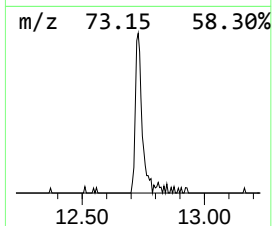
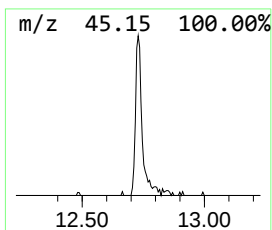
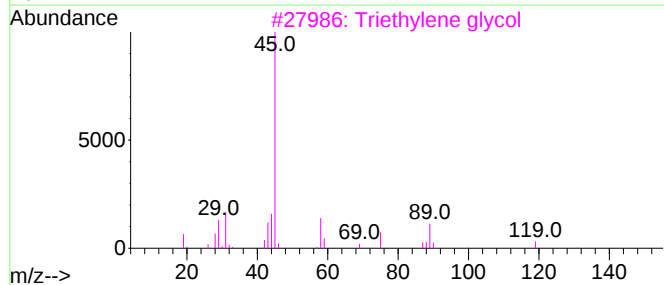
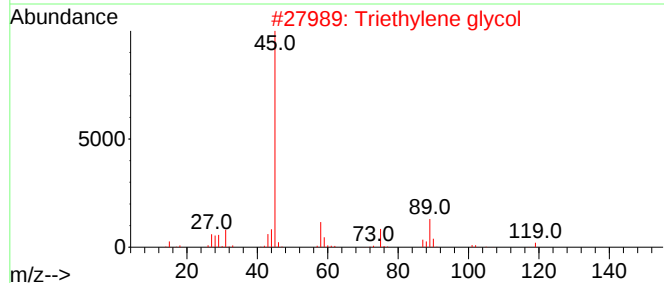
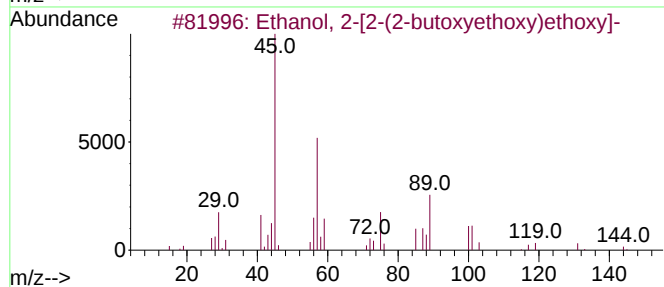
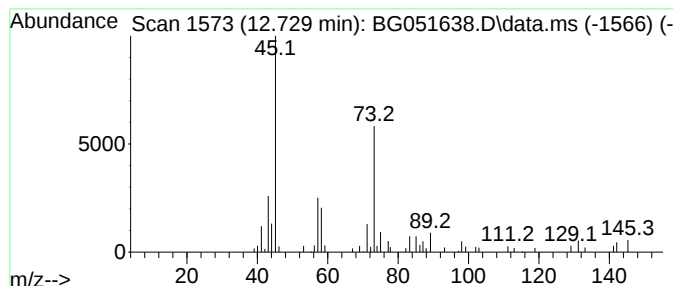
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Ethanol, 2-[2-(2-butoxyetho... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.729	3.92 ng/ul	62067	Naphthalene-d8	11.113

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-[2-(2-butoxyethoxy)et...	206	C10H22O4	000143-22-6	50
2			Triethylene glycol	150	C6H14O4	000112-27-6	50
3			Triethylene glycol	150	C6H14O4	000112-27-6	45
4			2-Pentanol, 3-chloro-4-methyl-, ...	136	C6H13ClO	074685-48-6	42
5			Triethylene glycol	150	C6H14O4	000112-27-6	40



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121721\
Data File : BG051638.D
Acq On : 18 Dec 2021 20:45
Operator : CG/JU
Sample : M5153-10
Misc :
ALS Vial : 54 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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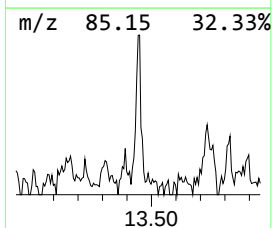
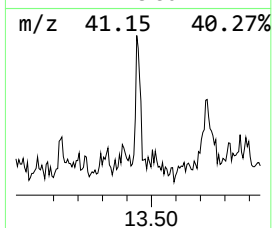
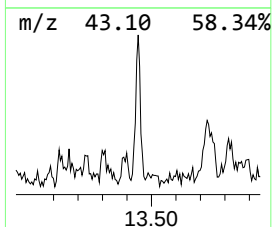
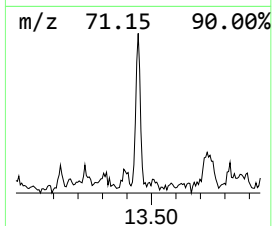
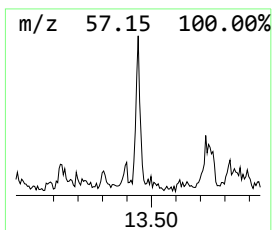
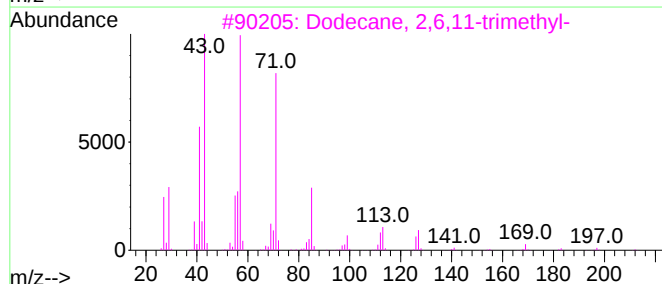
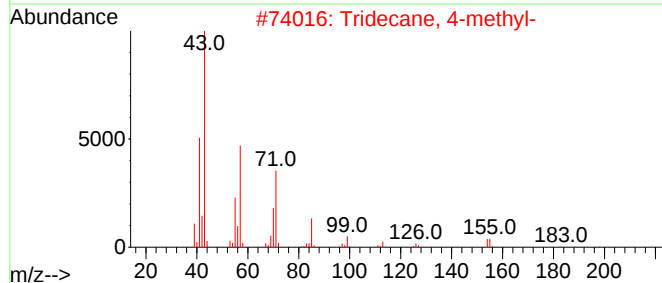
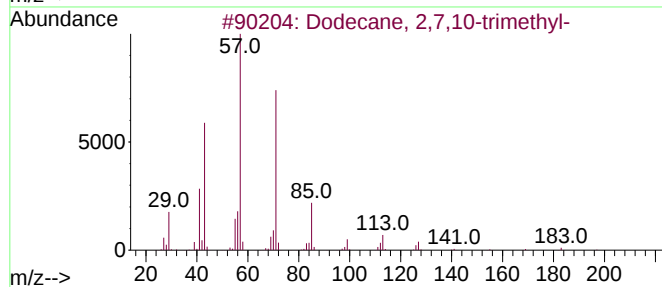
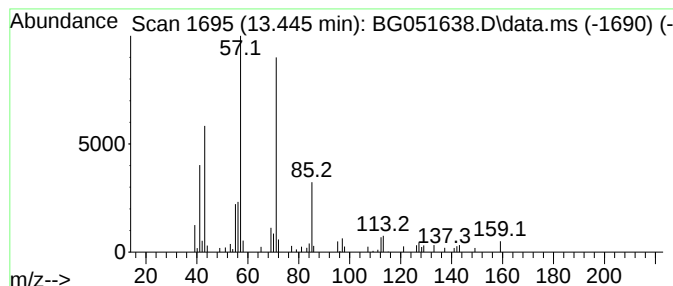
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.445	2.65 ng/ul	63936	Acenaphthene-d10	14.908

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	86
2			Tridecane, 4-methyl-	198	C14H30	026730-12-1	72
3			Dodecane, 2,6,11-trimethyl-	212	C15H32	031295-56-4	64
4			Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	58
5			Sulfurous acid, 2-ethylhexyl und...	348	C19H40O3S	1000309-19-4	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121721\
Data File : BG051638.D
Acq On : 18 Dec 2021 20:45
Operator : CG/JU
Sample : M5153-10
Misc :
ALS Vial : 54 Sample Multiplier: 1

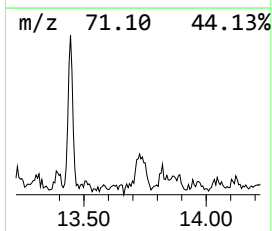
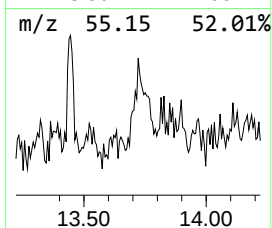
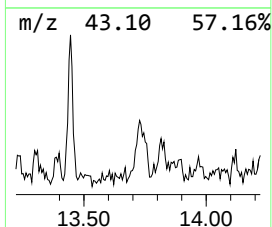
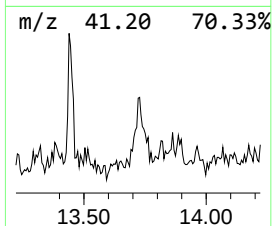
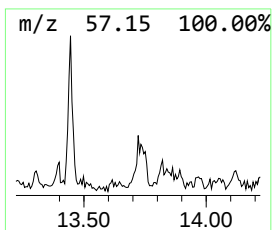
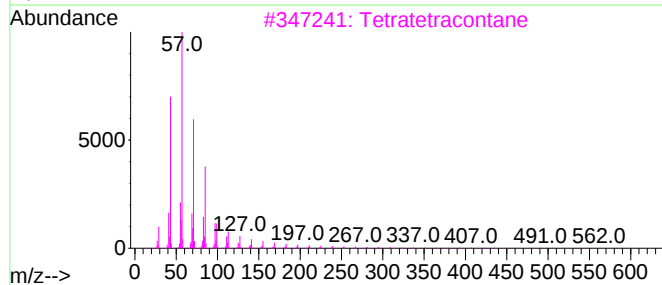
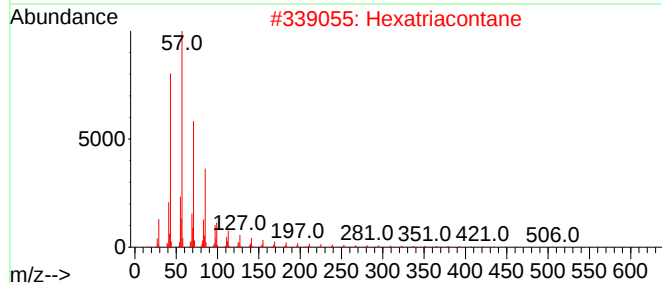
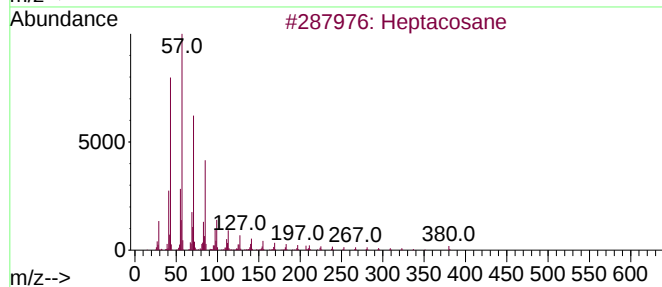
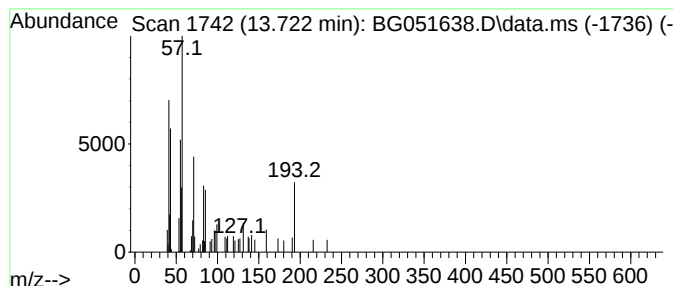
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121421.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.722	2.20 ng/ul	53115	Acenaphthene-d10	14.908	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptacosane	380	C27H56	000593-49-7	46
2	Hexatriacontane	507	C36H74	000630-06-8	43
3	Tetratetracontane	619	C44H90	007098-22-8	43
4	Octadecane, 1-bromo-	332	C18H37Br	000112-89-0	43
5	Hentriacontane	437	C31H64	000630-04-6	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121721\
Data File : BG051638.D
Acq On : 18 Dec 2021 20:45
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Sample : M5153-10
Misc :
ALS Vial : 54 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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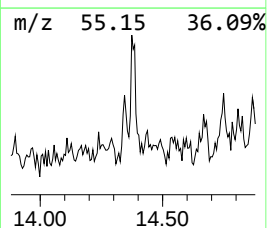
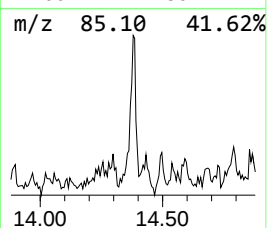
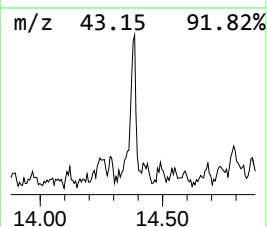
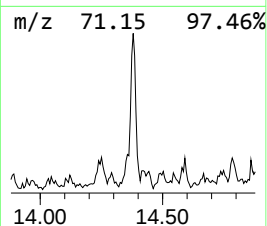
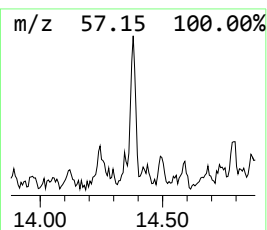
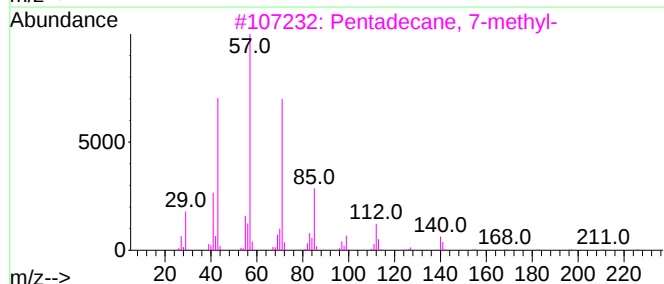
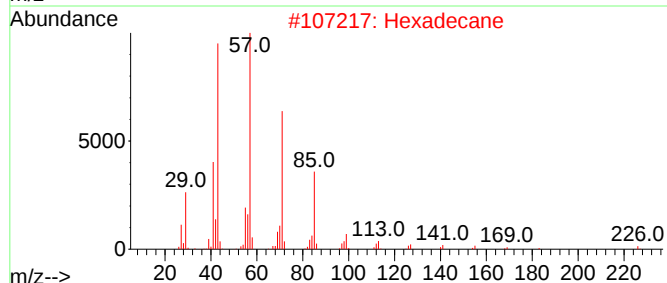
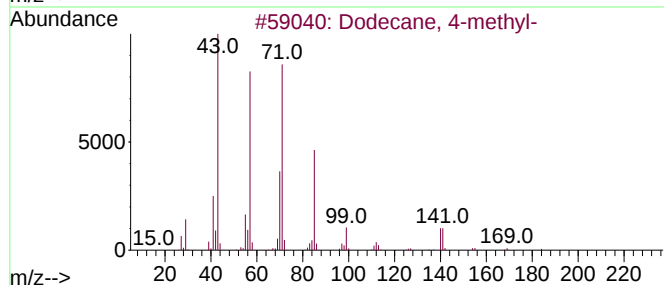
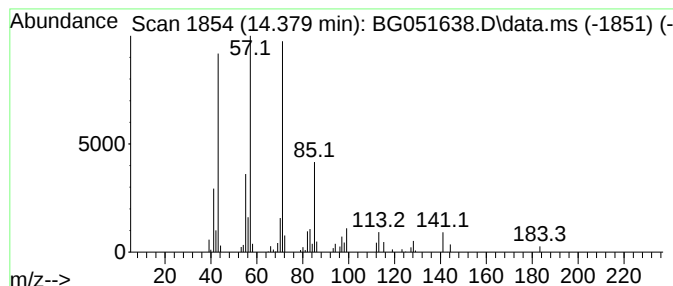
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.380	2.79 ng/ul	67347	Acenaphthene-d10	14.908

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecane, 4-methyl-	184	C13H28	006117-97-1	78
2			Hexadecane	226	C16H34	000544-76-3	72
3			Pentadecane, 7-methyl-	226	C16H34	006165-40-8	64
4			Sulfurous acid, decyl pentyl ester	292	C15H32O3S	1000309-14-3	64
5			Nonane, 5-(2-methylpropyl)-	184	C13H28	062185-53-9	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121721\
Data File : BG051638.D
Acq On : 18 Dec 2021 20:45
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Sample : M5153-10
Misc :
ALS Vial : 54 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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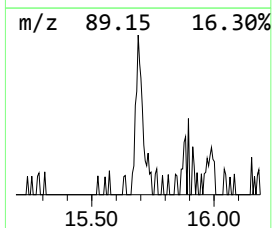
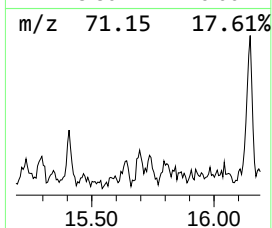
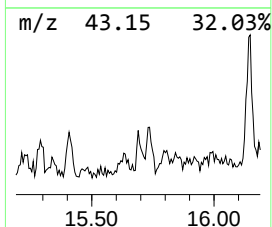
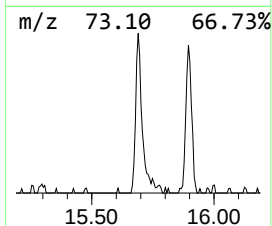
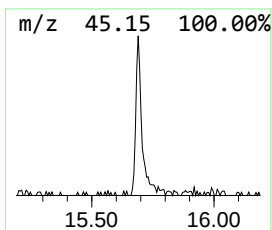
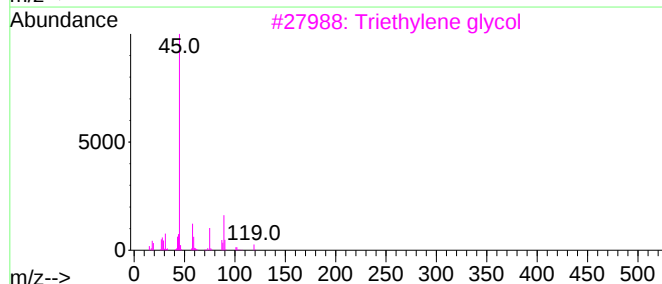
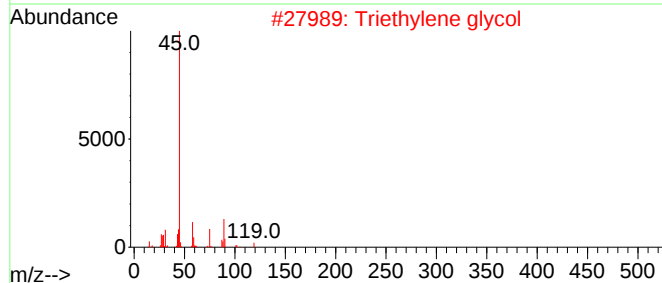
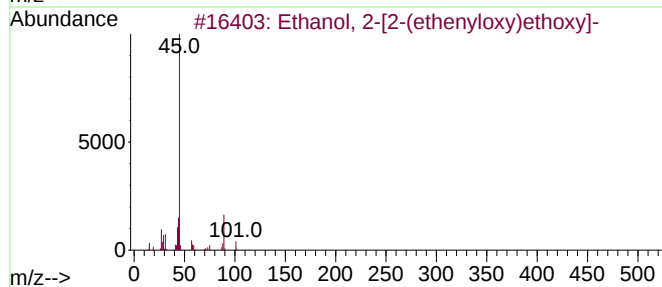
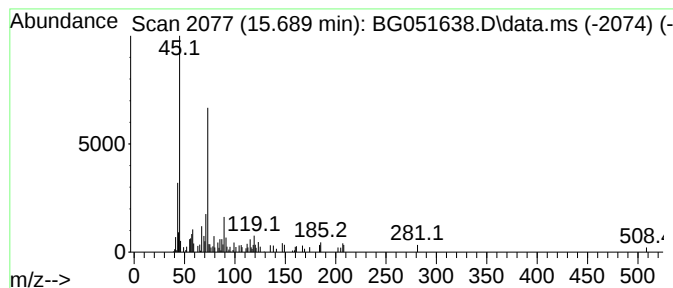
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Ethanol, 2-[2-(ethenyloxy)e... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.690	4.22 ng/ul	101773	Acenaphthene-d10	14.908

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-[2-(ethenyloxy)ethoxy]-	132	C6H12O3	000929-37-3	59
2			Triethylene glycol	150	C6H14O4	000112-27-6	53
3			Triethylene glycol	150	C6H14O4	000112-27-6	53
4			1-Pentanol, 5-methoxy-	118	C6H14O2	004799-62-6	50
5			Fumaric acid, bis(methoxymethyl)...	204	C8H12O6	1000197-25-6	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121721\
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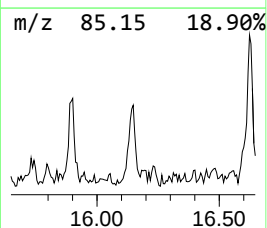
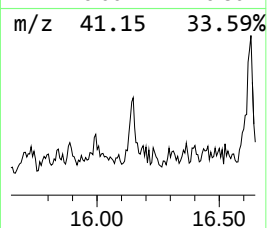
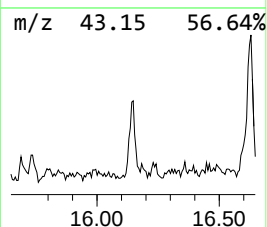
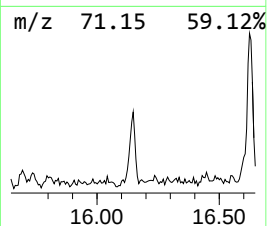
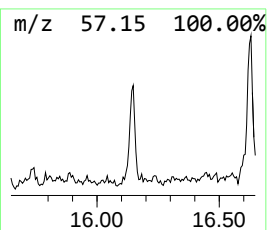
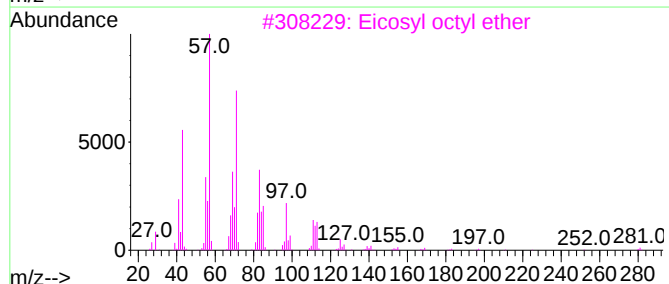
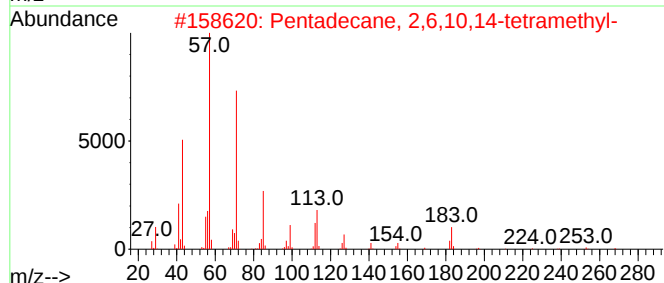
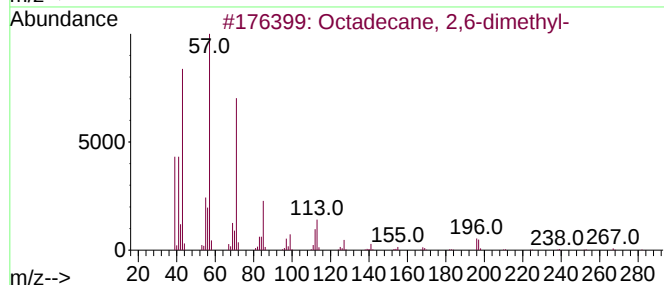
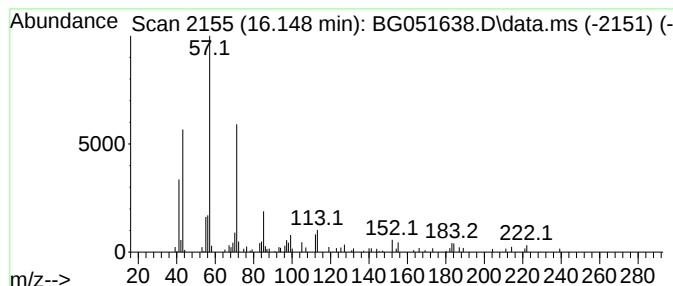
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121421.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
16.148	2.84 ng/ul	68536	Acenaphthene-d10	14.908	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecane, 2,6-dimethyl-	282	C20H42	075163-97-2	64
2	Pentadecane, 2,6,10,14-tetramethyl-	268	C19H40	001921-70-6	64
3	Eicosyl octyl ether	410	C28H58O	1000406-38-8	59
4	Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	59
5	Docosyl octyl ether	438	C30H62O	1000406-38-9	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121721\
Data File : BG051638.D
Acq On : 18 Dec 2021 20:45
Operator : CG/JU
Sample : M5153-10
Misc :
ALS Vial : 54 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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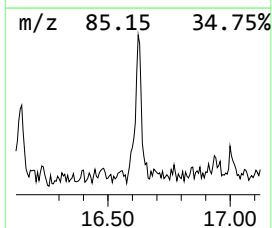
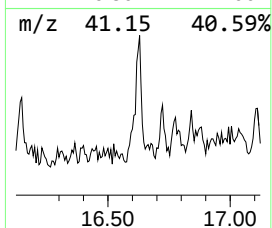
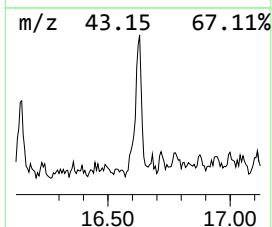
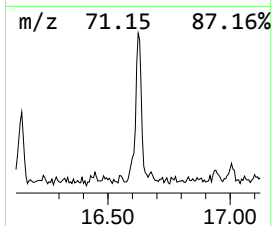
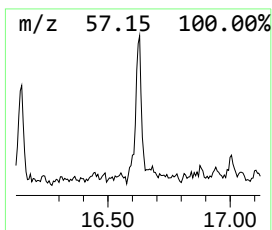
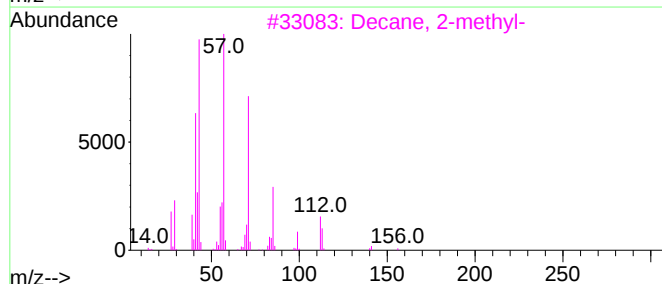
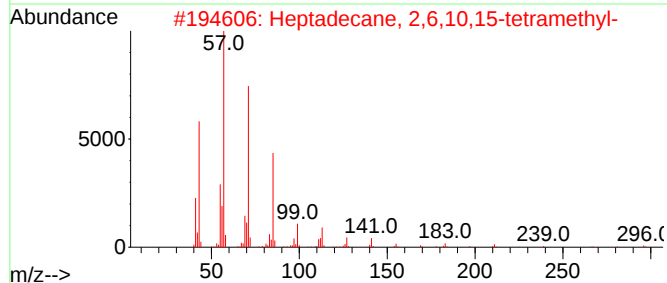
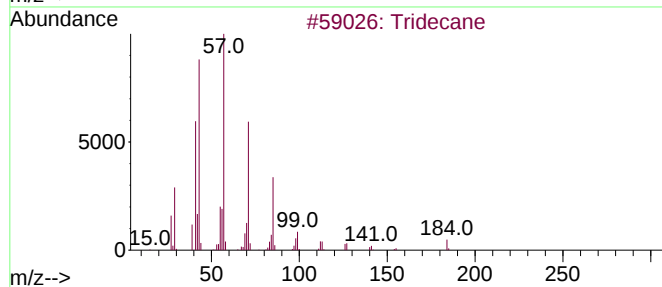
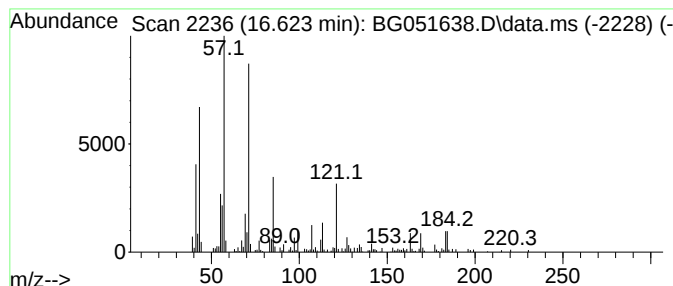
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.624	7.99 ng/ul	210388	Phenanthrene-d10	17.657

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tridecane	184	C13H28	000629-50-5	83
2		Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	80
3		Decane, 2-methyl-	156	C11H24	006975-98-0	74
4		Heptacosane	380	C27H56	000593-49-7	58
5		Nonane, 3-methyl-	142	C10H22	005911-04-6	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121721\
Data File : BG051638.D
Acq On : 18 Dec 2021 20:45
Operator : CG/JU
Sample : M5153-10
Misc :
ALS Vial : 54 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGL75

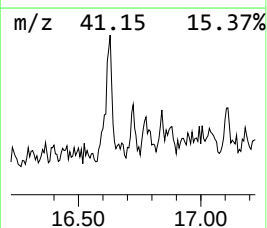
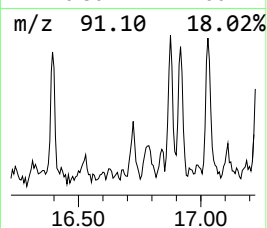
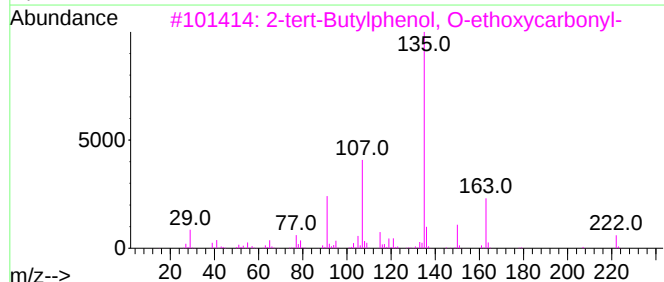
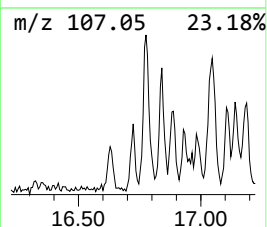
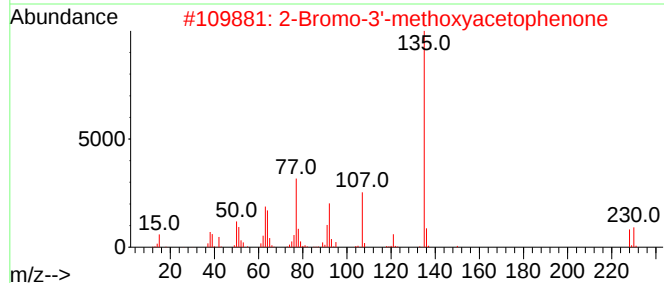
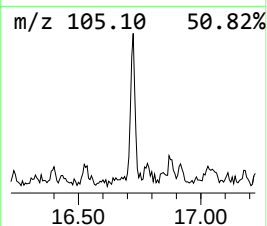
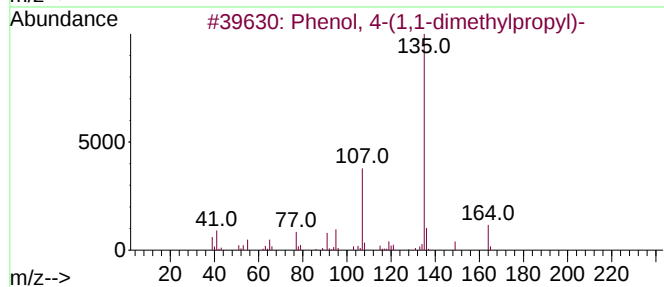
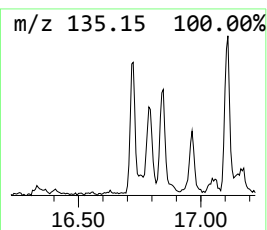
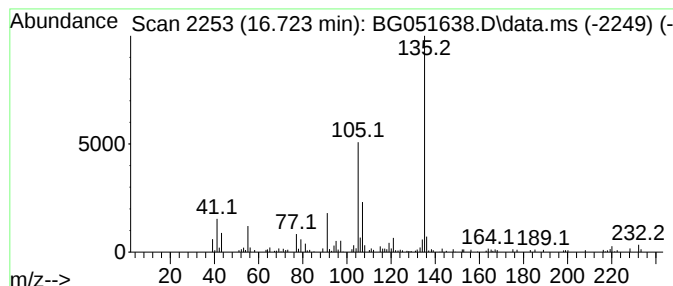
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 Phenol, 4-(1,1-dimethylprop... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.723	3.45 ng/ul	90861	Phenanthrene-d10	17.657

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	80
2			2-Bromo-3'-methoxyacetophenone	228	C9H9BrO2	005000-65-7	72
3			2-tert-Butylphenol, O-ethoxycarb...	222	C13H18O3	1000487-37-8	72
4			2-tert-Butylphenol, O-(n-propylo...	236	C14H20O3	1000487-37-9	72
5			Benzeneacetonitrile, 3-fluoro-	135	C8H6FN	000501-00-8	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121721\
Data File : BG051638.D
Acq On : 18 Dec 2021 20:45
Operator : CG/JU
Sample : M5153-10
Misc :
ALS Vial : 54 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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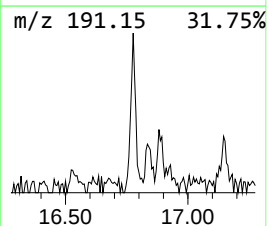
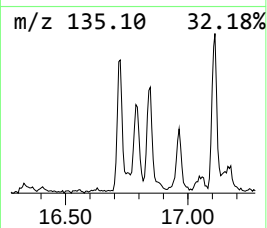
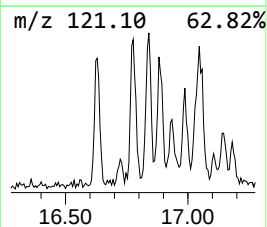
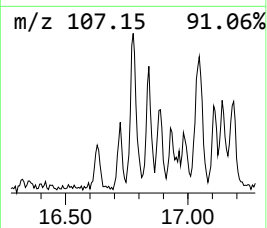
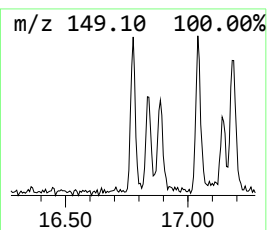
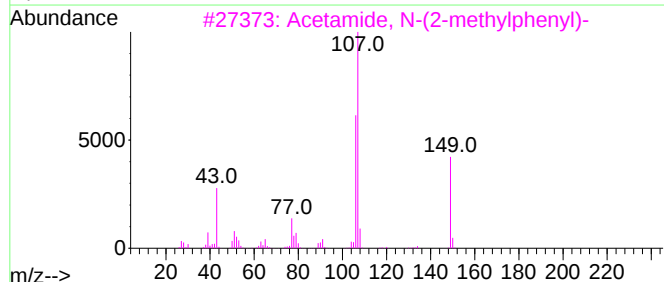
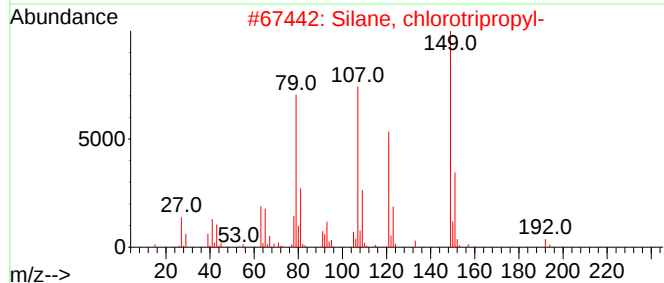
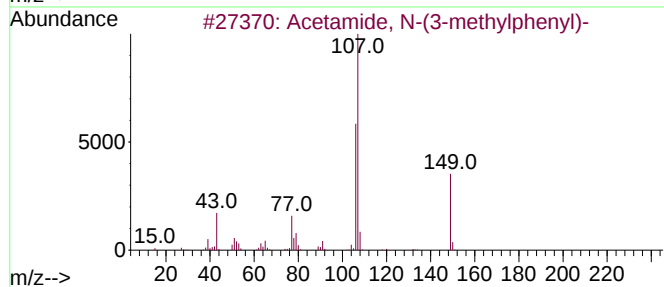
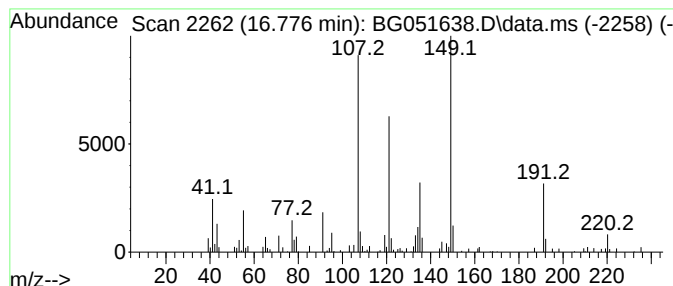
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 Acetamide, N-(3-methylphenyl)- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.776	4.76 ng/ul	125320	Phenanthrene-d10	17.657

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acetamide, N-(3-methylphenyl)-	149	C9H11NO	000537-92-8	62
2			Silane, chlorotripropyl-	192	C9H21ClSi	000995-25-5	59
3			Acetamide, N-(2-methylphenyl)-	149	C9H11NO	000120-66-1	58
4			Acetamide, N-(2-methylphenyl)-	149	C9H11NO	000120-66-1	46
5			Acetamide, N-(4-methylphenyl)-	149	C9H11NO	000103-89-9	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121721\
Data File : BG051638.D
Acq On : 18 Dec 2021 20:45
Operator : CG/JU
Sample : M5153-10
Misc :
ALS Vial : 54 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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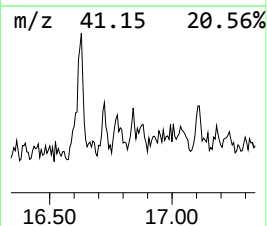
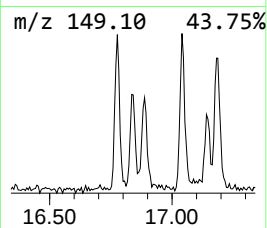
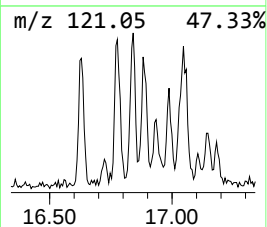
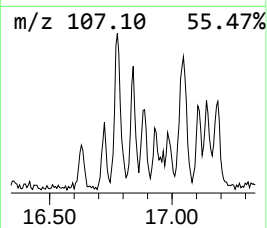
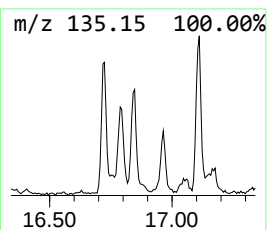
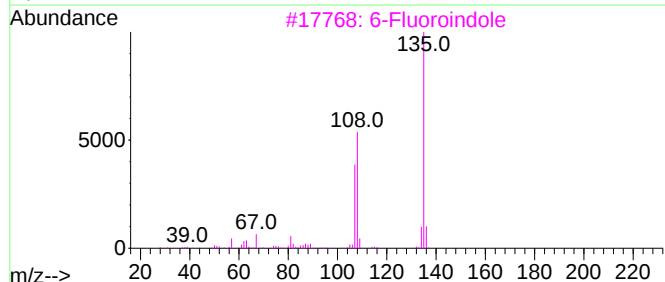
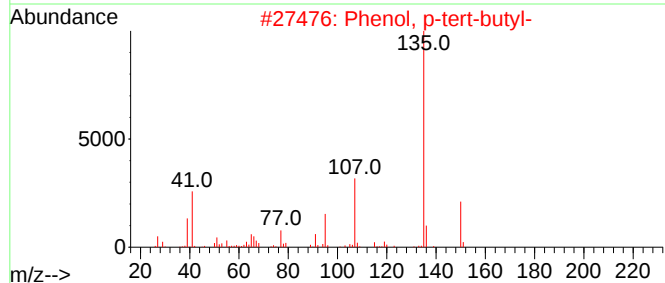
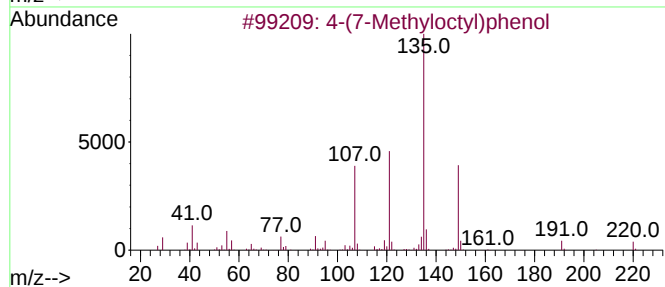
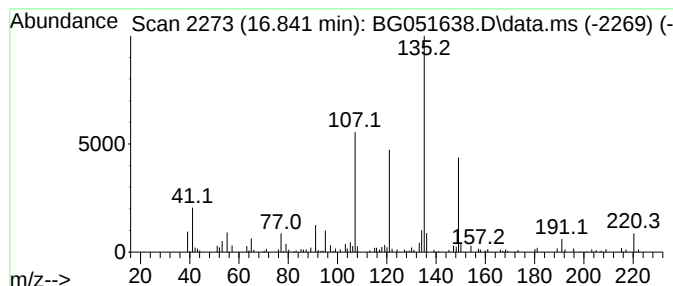
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 4-(7-Methyloctyl)phenol Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.841	2.94 ng/ul	77414	Phenanthrene-d10	17.657

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		4-(7-Methyloctyl)phenol	220	C15H24O	024518-48-7	64
2		Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	53
3		6-Fluoroindole	135	C8H6FN	000399-51-9	50
4		Benzeneacetonitrile, 3-fluoro-	135	C8H6FN	000501-00-8	47
5		Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121721\
Data File : BG051638.D
Acq On : 18 Dec 2021 20:45
Operator : CG/JU
Sample : M5153-10
Misc :
ALS Vial : 54 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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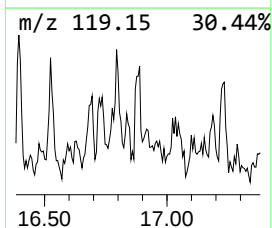
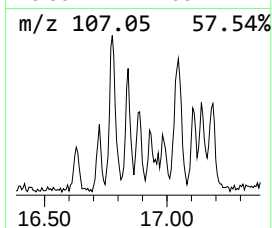
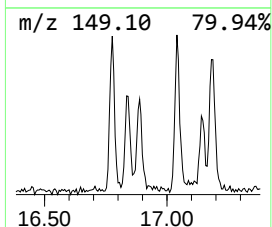
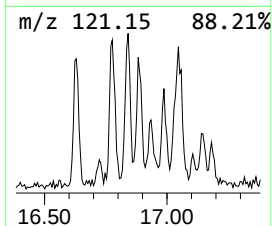
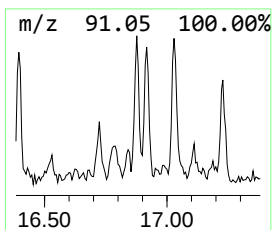
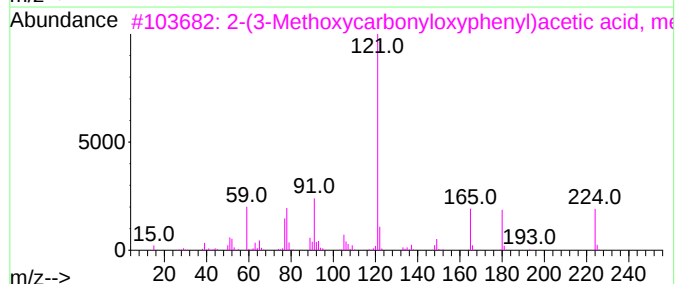
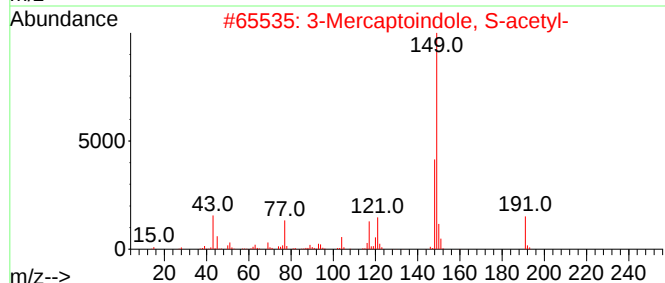
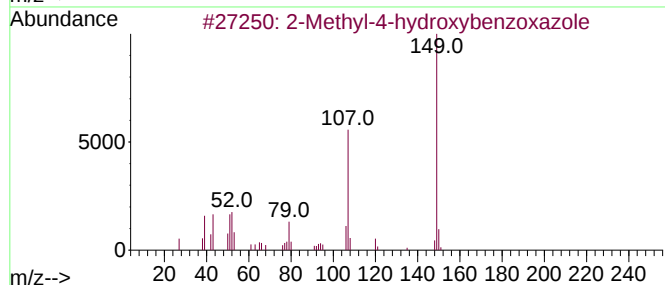
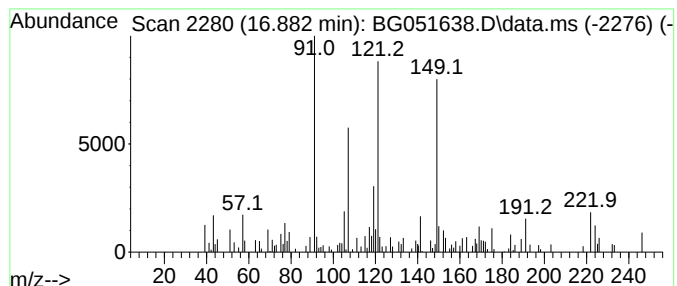
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 unknown-02 Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.882	2.50 ng/ul	65699	Phenanthrene-d10	17.657

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Methyl-4-hydroxybenzoxazole	149	C8H7NO2	051110-60-2	35
2			3-Mercaptoindole, S-acetyl-	191	C10H9NOS	1081542-19-9	25
3			2-(3-Methoxycarbonyloxyphenyl)ac...	224	C11H12O5	1000487-68-3	25
4			Benzene, 1-isocyanato-3-methoxy-	149	C8H7NO2	018908-07-1	22
5			Acetamide, N-(4-methylphenyl)-	149	C9H11NO	000103-89-9	22



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121721\
Data File : BG051638.D
Acq On : 18 Dec 2021 20:45
Operator : CG/JU
Sample : M5153-10
Misc :
ALS Vial : 54 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121421.M
Quant Title : SVOA CALIBRATION

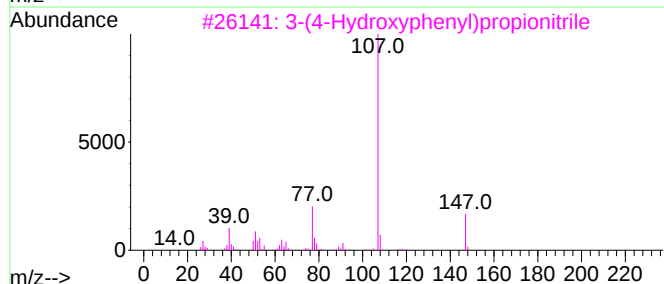
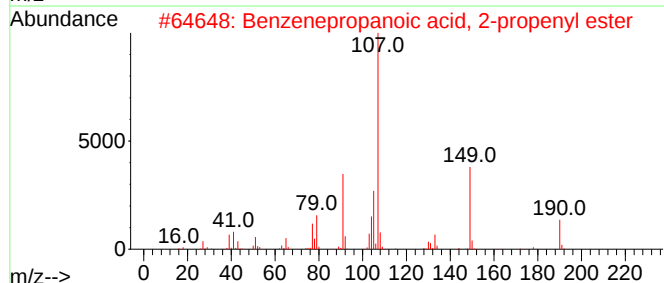
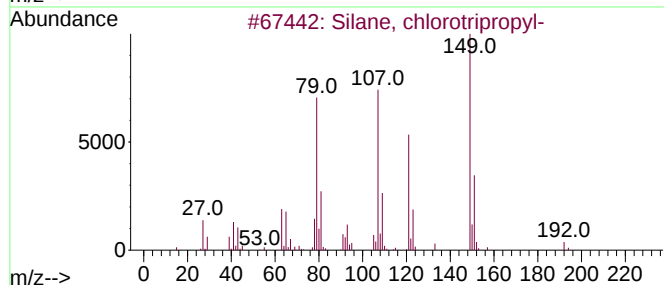
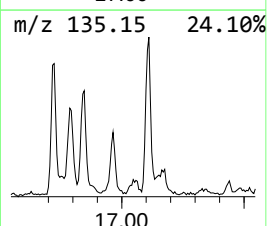
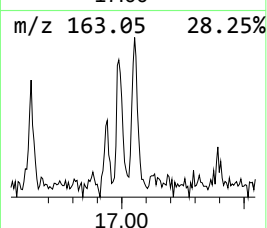
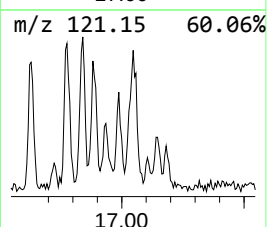
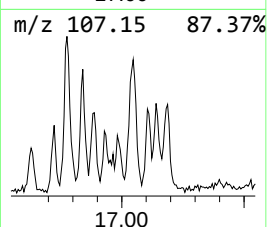
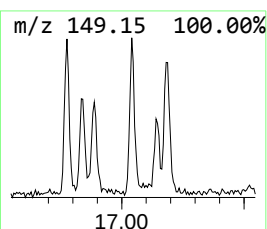
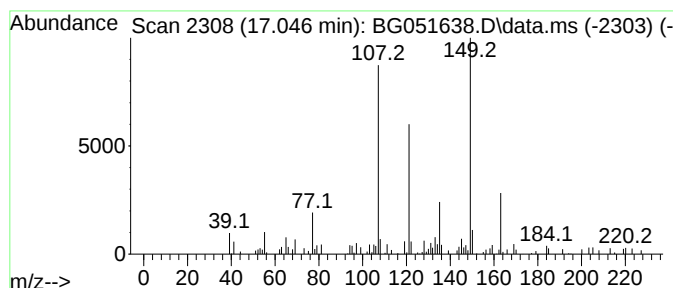
TIC Library : C:\DATABASE\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 15 Silane, chlorotripropyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.047	4.59 ng/ul	120710	Phenanthrene-d10	17.657

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Silane, chlorotripropyl-	192	C9H21ClSi	000995-25-5	53
2			Benzenepropanoic acid, 2-propeny...	190	C12H14O2	015814-45-6	38
3			3-(4-Hydroxyphenyl)propionitrile	147	C9H9NO	017362-17-3	27
4			Phenol, 2-(1,1-dimethylethyl)-4-...	164	C11H16O	002409-55-4	27
5			Pyridine, pentamethyl-	149	C10H15N	003748-83-2	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121721\
Data File : BG051638.D
Acq On : 18 Dec 2021 20:45
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Sample : M5153-10
Misc :
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ClientSampleId :
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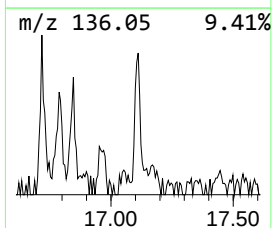
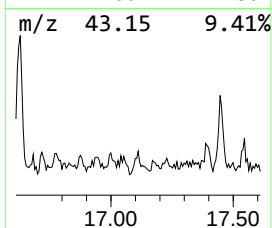
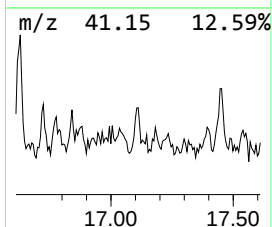
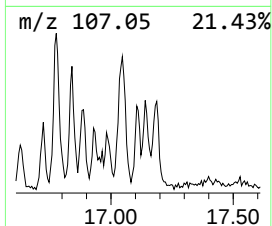
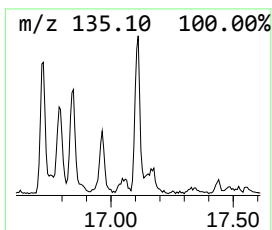
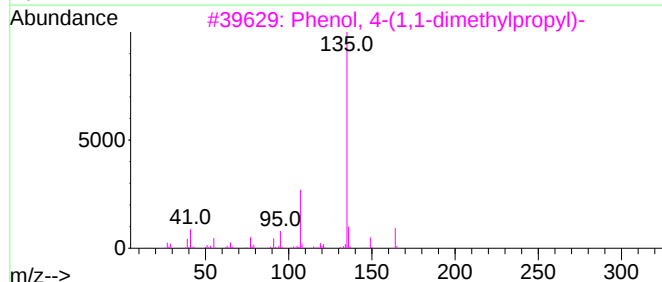
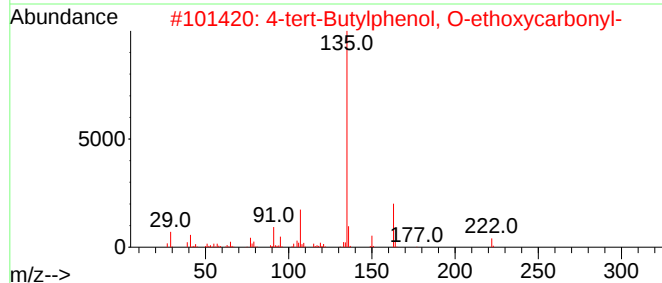
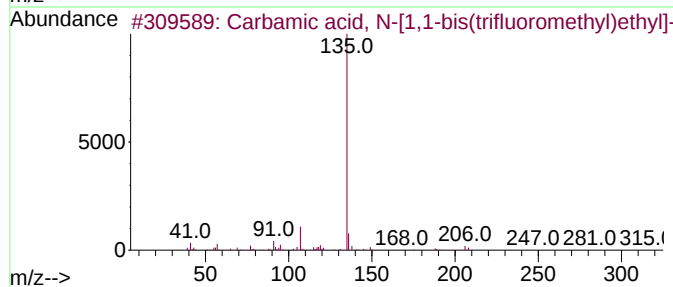
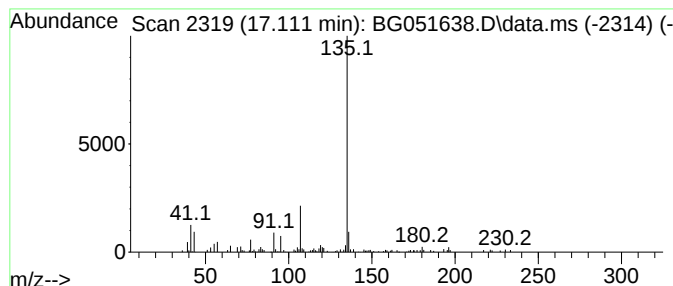
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 16 Carbamic acid, N-[1,1-bis(t... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.111	2.83 ng/ul	74602	Phenanthrene-d10	17.657

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Carbamic acid, N-[1,1-bis(triflu...	413	C19H25F6N02	296242-69-8	78
2			4-tert-Butylphenol, O-ethoxycarb...	222	C13H18O3	026177-66-2	78
3			Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	72
4			Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	72
5			Phenol, 2-(1,1,3,3-tetramethylbu...	206	C14H22O	003884-95-5	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121721\
Data File : BG051638.D
Acq On : 18 Dec 2021 20:45
Operator : CG/JU
Sample : M5153-10
Misc :
ALS Vial : 54 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGL75

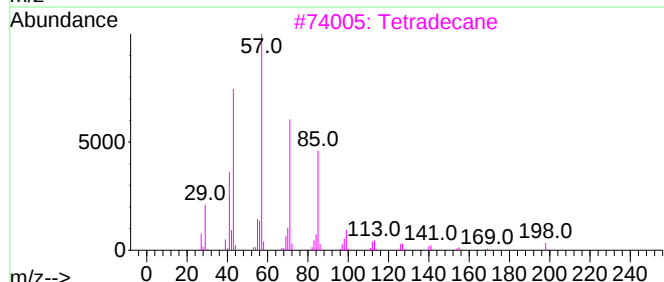
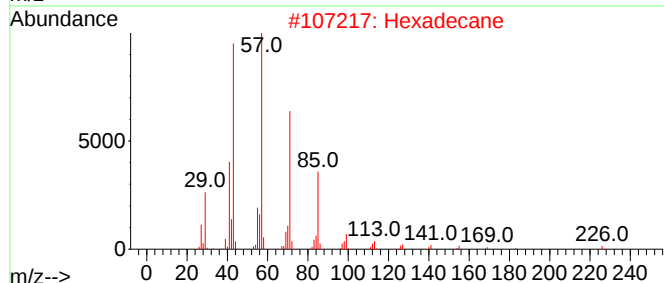
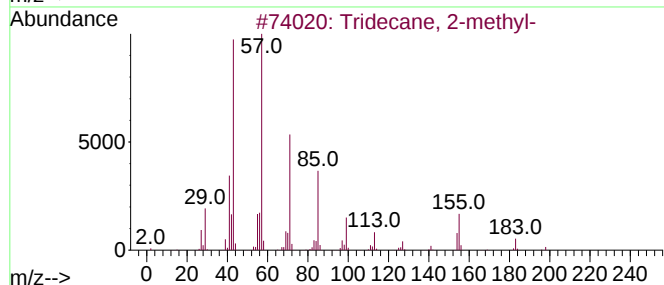
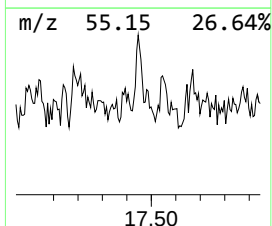
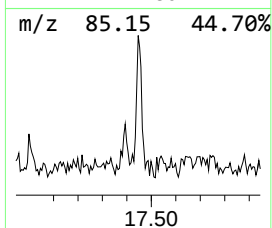
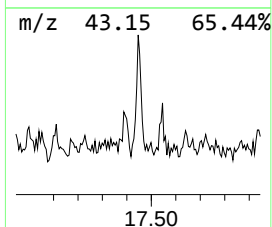
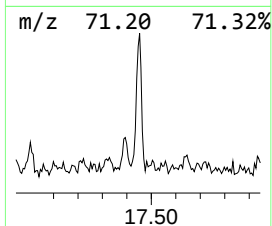
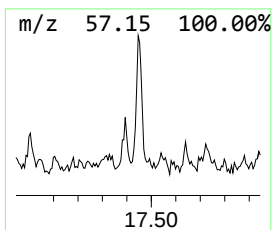
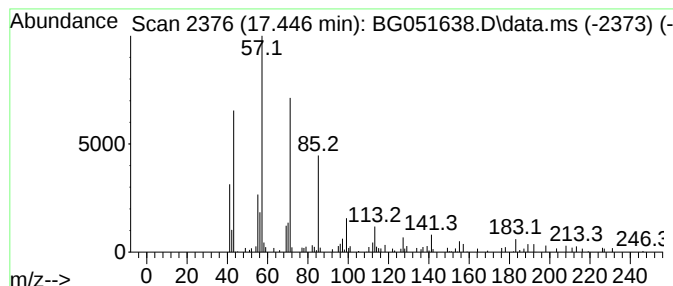
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 17 (DEL) Alkane: Straight-Chai... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.446	2.77 ng/ul	72814	Phenanthrene-d10	17.657

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tridecane, 2-methyl-	198	C14H30	001560-96-9	90
2		Hexadecane	226	C16H34	000544-76-3	87
3		Tetradecane	198	C14H30	000629-59-4	87
4		3-Ethyl-2,6,10-trimethylundecane	226	C16H34	1000432-25-9	83
5		Tetradecane	198	C14H30	000629-59-4	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121721\
Data File : BG051638.D
Acq On : 18 Dec 2021 20:45
Operator : CG/JU
Sample : M5153-10
Misc :
ALS Vial : 54 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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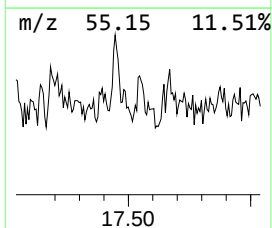
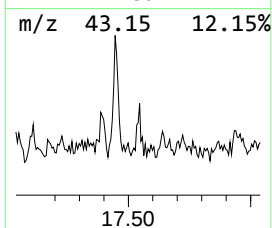
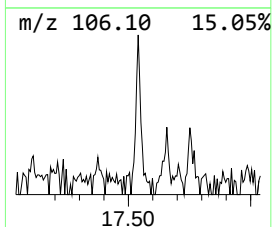
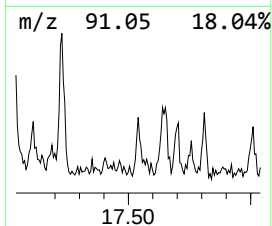
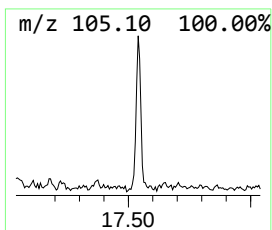
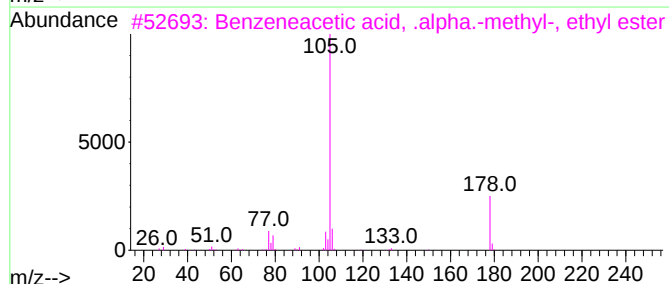
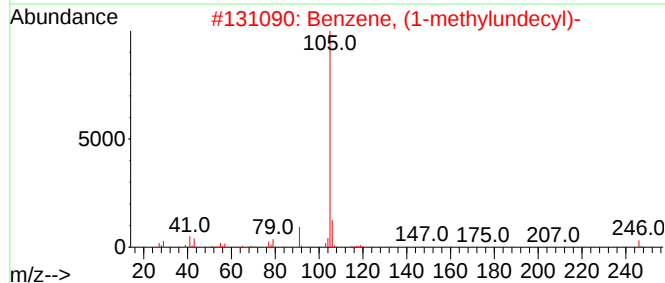
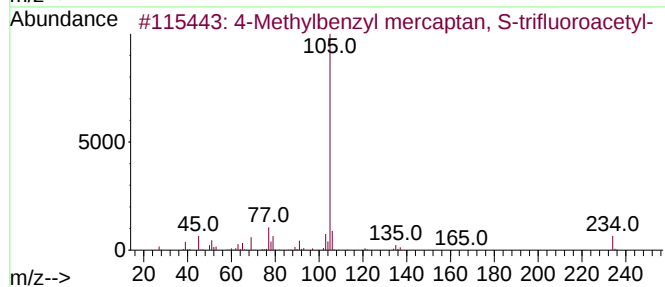
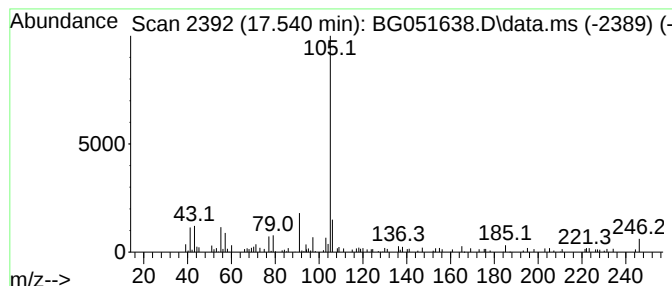
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 18 4-Methylbenzyl mercaptan, S... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.540	2.46 ng/ul	64632	Phenanthrene-d10	17.657

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-Methylbenzyl mercaptan, S-trif...	234	C10H9F3OS	1000470-56-8	53
2			Benzene, (1-methylundecyl)-	246	C18H30	002719-61-1	53
3			Benzeneacetic acid, .alpha.-meth...	178	C11H14O2	002510-99-8	53
4			Methyl N-(4-methylbenzyl)-N-(vin...	313	C14H19NO5S	1000481-55-4	53
5			Ethyl meta-tolylacetate	178	C11H14O2	040061-55-0	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121721\
Data File : BG051638.D
Acq On : 18 Dec 2021 20:45
Operator : CG/JU
Sample : M5153-10
Misc :
ALS Vial : 54 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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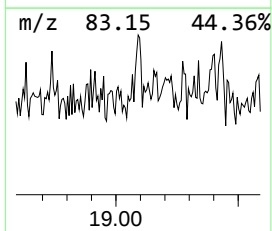
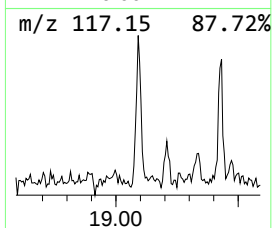
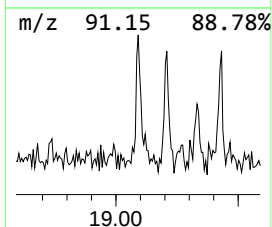
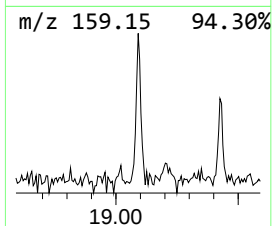
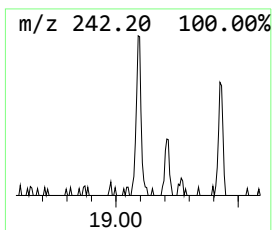
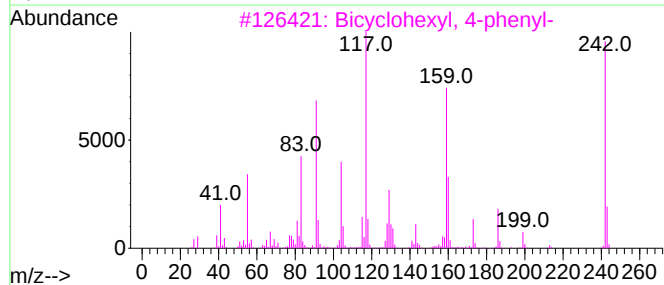
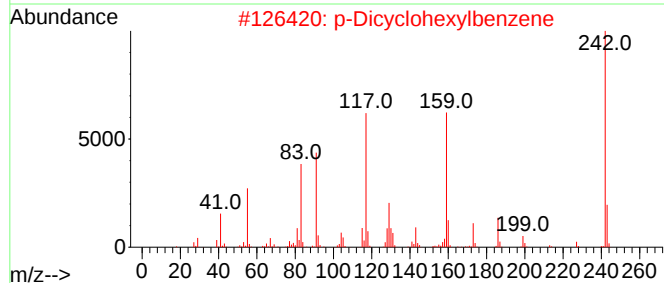
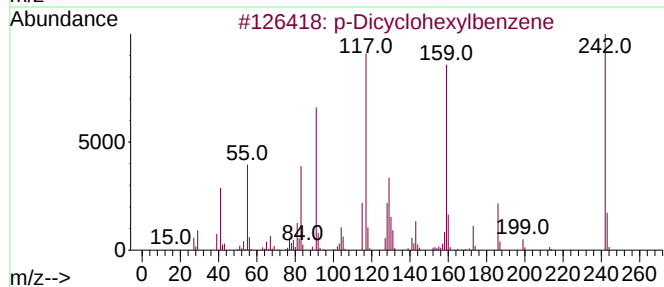
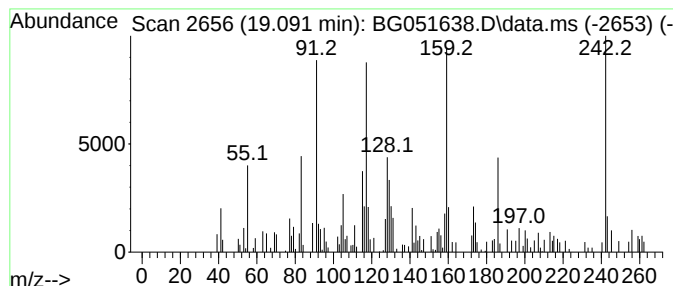
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 19 p-Dicyclohexylbenzene Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.091	2.03 ng/ul	53485	Phenanthrene-d10	17.657

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			p-Dicyclohexylbenzene	242	C18H26	001087-02-1	93
2			p-Dicyclohexylbenzene	242	C18H26	001087-02-1	86
3			Bicyclohexyl, 4-phenyl-	242	C18H26	020273-27-2	58
4			p-Dicyclohexylbenzene	242	C18H26	001087-02-1	53
5			1(2H)-Isoquinolinone, 2-methyl-	159	C10H9NO	004594-71-2	18



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121721\
Data File : BG051638.D
Acq On : 18 Dec 2021 20:45
Operator : CG/JU
Sample : M5153-10
Misc :
ALS Vial : 54 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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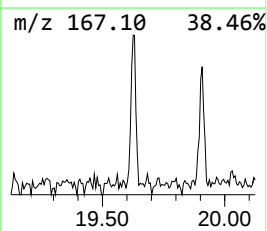
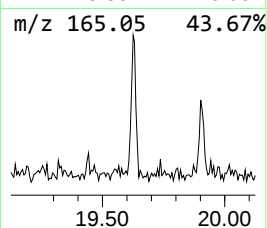
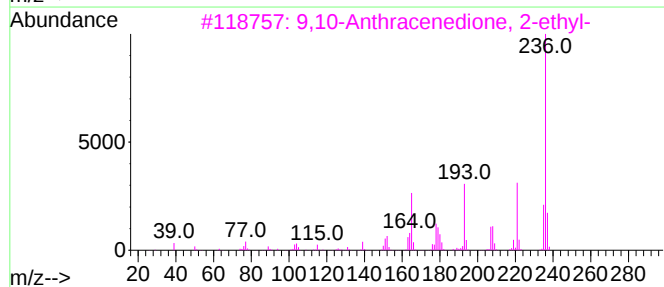
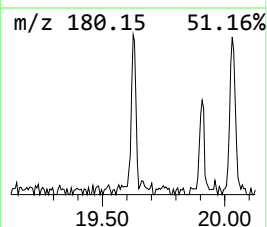
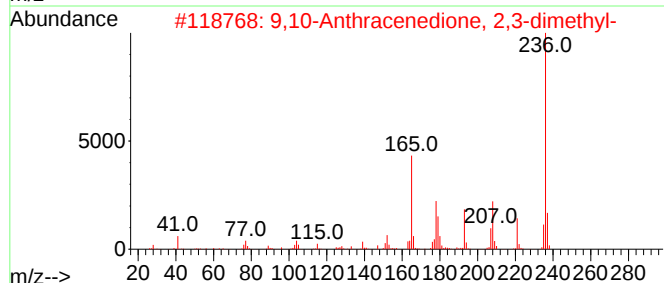
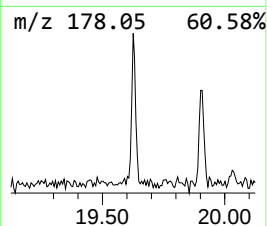
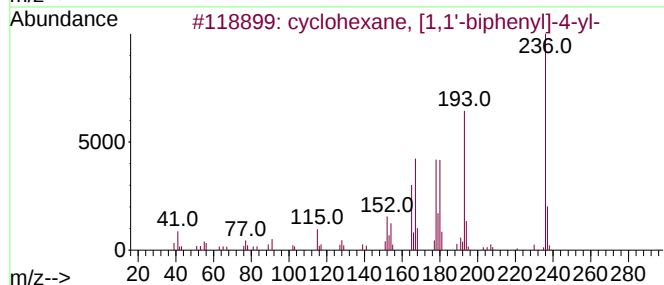
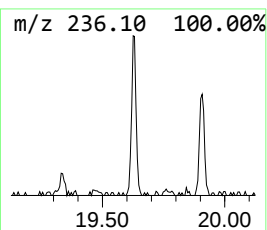
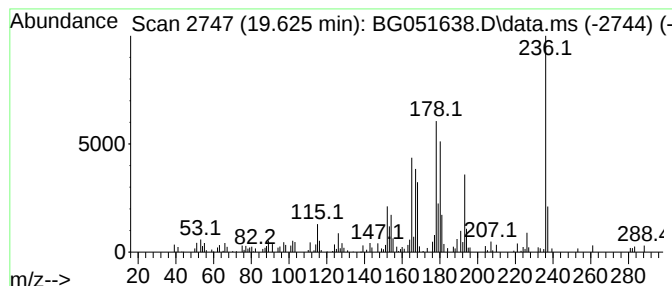
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 20 cyclohexane, [1,1'-biphenyl]... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.625	3.49 ng/ul	91836	Phenanthrene-d10	17.657

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			cyclohexane, [1,1'-biphenyl]-4-yl-	236	C18H20	1000401-12-4	76
2			9,10-Anthracenedione, 2,3-dimethyl-	236	C16H12O2	006531-35-7	49
3			9,10-Anthracenedione, 2-ethyl-	236	C16H12O2	000084-51-5	43
4			7-Trimethylsilyloxychromanone	236	C12H16O3Si	1000451-97-2	43
5			9,10-Anthracenedione, 1,4-dimethyl-	236	C16H12O2	001519-36-4	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121721\
Data File : BG051638.D
Acq On : 18 Dec 2021 20:45
Operator : CG/JU
Sample : M5153-10
Misc :
ALS Vial : 54 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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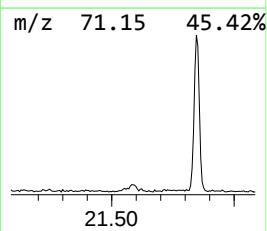
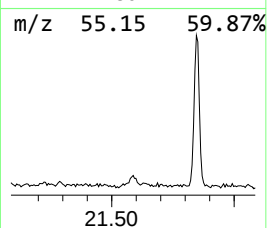
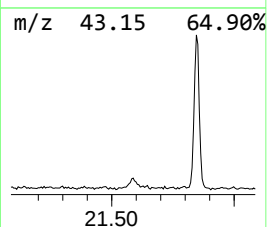
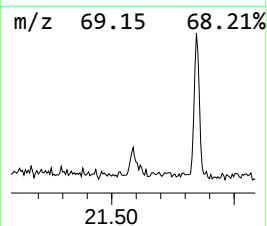
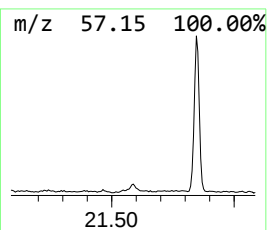
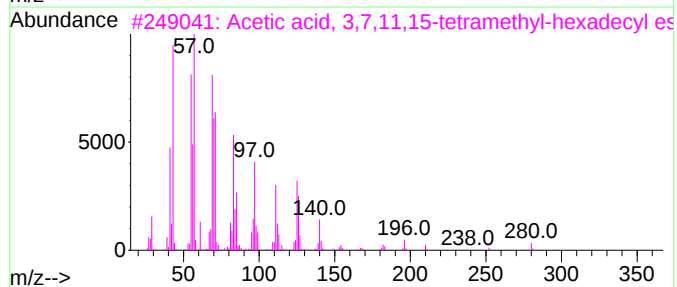
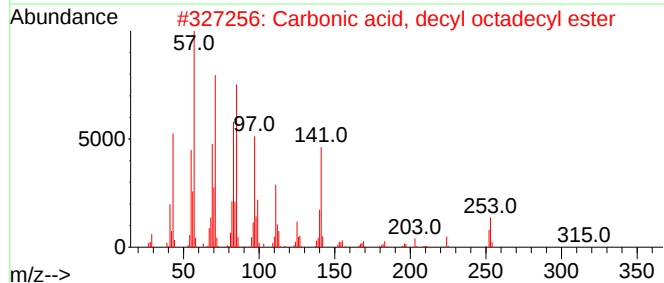
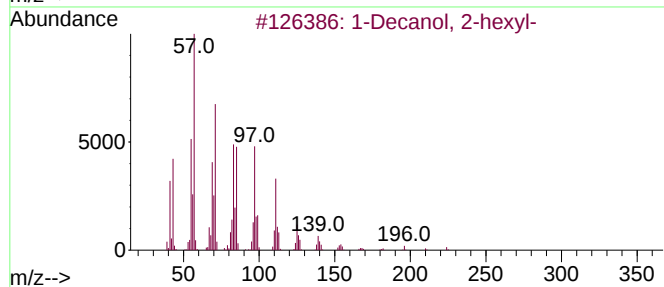
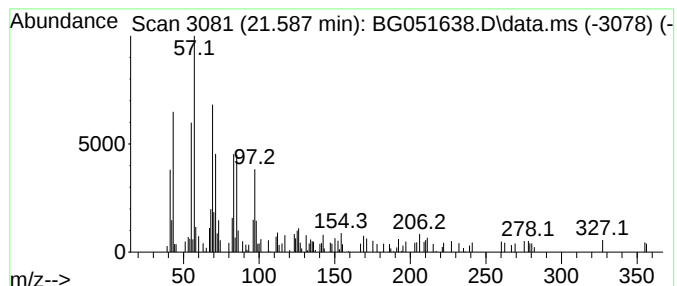
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 21 1-Decanol, 2-hexyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.587	3.01 ng/ul	88930	Chrysene-d12	21.975

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Decanol, 2-hexyl-			242	C16H34O	002425-77-6	80
2	Carbonic acid, decyl octadecyl e...			454	C29H58O3	1000383-16-7	80
3	Acetic acid, 3,7,11,15-tetrameth...			340	C22H44O2	1000193-63-0	58
4	Trichloroacetic acid, tridecyl e...			344	C15H27Cl3O2	074339-51-8	49
5	Heptadecyl heptafluorobutyrate			452	C21H35F7O2	959085-66-6	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121721\
Data File : BG051638.D
Acq On : 18 Dec 2021 20:45
Operator : CG/JU
Sample : M5153-10
Misc :
ALS Vial : 54 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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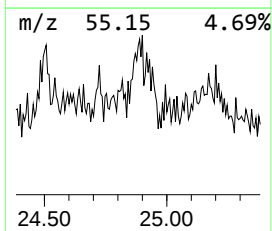
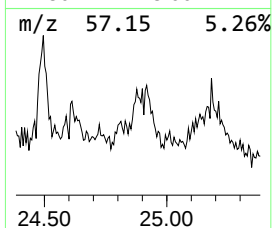
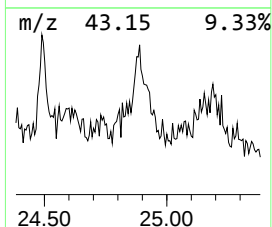
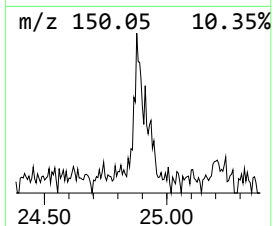
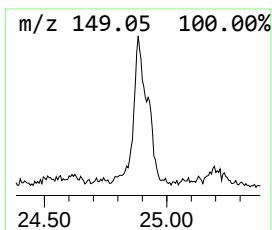
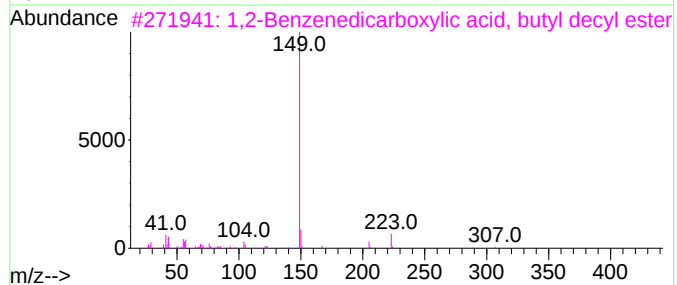
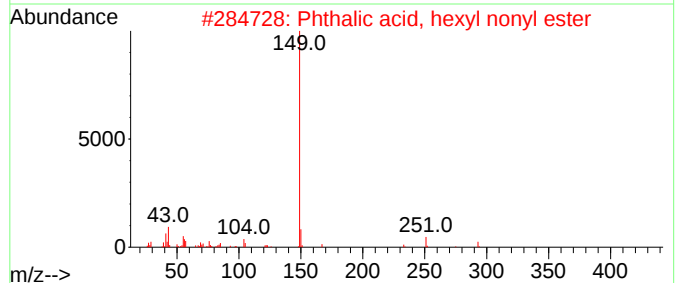
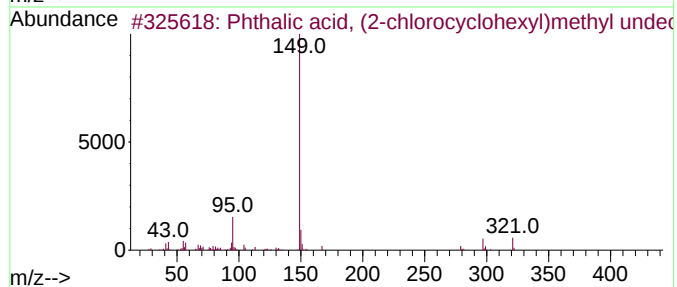
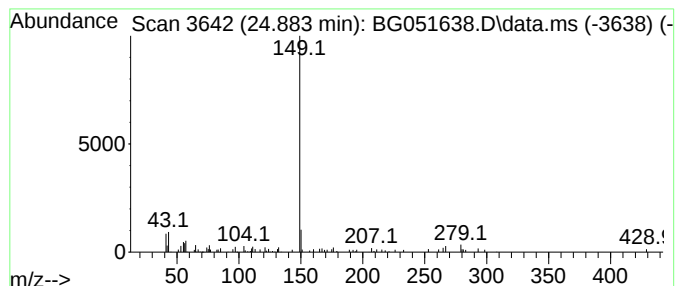
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 22 Phthalic acid, (2-chlorocyclohexyl)methyl ester Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.883	2.17 ng/ul	61349	Perylene-d12	25.476

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phthalic acid, (2-chlorocyclohexyl)methyl ester	450	C26H39ClO4	1000315-34-9	72
2			Phthalic acid, hexyl nonyl ester	376	C23H36O4	1000309-01-1	64
3			1,2-Benzenedicarboxylic acid, butyl decyl ester	362	C22H34O4	000089-19-0	64
4			Phthalic acid, 6-ethyl-3-octyl hexyl ester	404	C25H40O4	1000315-17-7	64
5			Phthalic acid, 2-chloropropyl tetradecyl ester	438	C25H39ClO4	1000356-83-5	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121721\
 Data File : BG051638.D
 Acq On : 18 Dec 2021 20:45
 Operator : CG/JU
 Sample : M5153-10
 Misc :
 ALS Vial : 54 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121421.M
 Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST0.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown-01	8.470	4.1	ng/ul	53136	1	8.276	258809	20.0
Sulfurous acid,...	8.529	2.0	ng/ul	26199	1	8.276	258809	20.0
(DEL) Alkane: S...	12.124	3.6	ng/ul	56549	2	11.113	316486	20.0
Ethanol, 2-[2-(...	12.729	3.9	ng/ul	62067	2	11.113	316486	20.0
(DEL) Alkane: S...	13.445	2.6	ng/ul	63936	3	14.908	482786	20.0
(DEL) Alkane: S...	13.722	2.2	ng/ul	53115	3	14.908	482786	20.0
(DEL) Alkane: S...	14.380	2.8	ng/ul	67347	3	14.908	482786	20.0
Ethanol, 2-[2-(...	15.690	4.2	ng/ul	101773	3	14.908	482786	20.0
(DEL) Alkane: S...	16.148	2.8	ng/ul	68536	3	14.908	482786	20.0
(DEL) Alkane: S...	16.624	8.0	ng/ul	210388	4	17.657	526370	20.0
Phenol, 4-(1,1-...	16.723	3.5	ng/ul	90861	4	17.657	526370	20.0
Acetamide, N-(3...	16.776	4.8	ng/ul	125320	4	17.657	526370	20.0
4-(7-Methylocty...	16.841	2.9	ng/ul	77414	4	17.657	526370	20.0
unknown-02	16.882	2.5	ng/ul	65699	4	17.657	526370	20.0
Silane, chlorot...	17.047	4.6	ng/ul	120710	4	17.657	526370	20.0
Carbamic acid, ...	17.111	2.8	ng/ul	74602	4	17.657	526370	20.0
(DEL) Alkane: S...	17.446	2.8	ng/ul	72814	4	17.657	526370	20.0
4-Methylbenzyl ...	17.540	2.5	ng/ul	64632	4	17.657	526370	20.0
p-Dicyclohexylb...	19.091	2.0	ng/ul	53485	4	17.657	526370	20.0
cyclohexane, [1...	19.625	3.5	ng/ul	91836	4	17.657	526370	20.0
1-Decanol, 2-he...	21.587	3.0	ng/ul	88930	5	21.975	590250	20.0
Phthalic acid, ...	24.883	2.2	ng/ul	61349	6	25.476	566089	20.0