

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG112321\
 Data File : BG051225.D
 Acq On : 25 Nov 2021 00:21
 Operator : CG/JU
 Sample : M4725-01
 Misc :
 ALS Vial : 47 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 F4L00

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-BG112321.M
 Title : SVOA CALIBRATION

Signal : TIC: BG051226.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.541	4	9	16	rBV2	6981	11010	1.12%	0.097%
2	3.976	78	83	91	rBV	47232	83507	8.51%	0.735%
3	4.311	134	140	146	rBV4	2565	4032	0.41%	0.035%
4	6.808	560	565	569	rVB4	3197	4676	0.48%	0.041%
5	7.360	652	659	673	rVB3	32097	73002	7.44%	0.643%
6	7.513	676	685	695	rBV	113448	196009	19.97%	1.725%
7	7.730	715	722	734	rBV	95672	167938	17.11%	1.478%
8	7.960	753	761	768	rVB4	1982	4185	0.43%	0.037%
9	8.200	794	802	813	rBV	115090	207508	21.14%	1.827%
10	8.911	915	923	930	rBV	81276	151682	15.46%	1.335%
11	8.964	930	932	938	rVV	10176	13543	1.38%	0.119%
12	9.376	995	1002	1015	rBV	148157	271220	27.64%	2.388%
13	9.511	1021	1025	1032	rVB4	2778	6223	0.63%	0.055%
14	9.716	1055	1060	1065	rVB2	5559	7974	0.81%	0.070%
15	10.098	1119	1125	1134	rBV	95655	180489	18.39%	1.589%
16	10.651	1212	1219	1228	rBV	139289	252970	25.78%	2.227%
17	10.774	1234	1240	1252	rBV3	29624	75344	7.68%	0.663%
18	11.027	1275	1283	1296	rBV	185640	342255	34.87%	3.013%
19	11.168	1299	1307	1321	rBV	131377	267094	27.22%	2.351%
20	11.291	1322	1328	1337	rVB	34864	59706	6.08%	0.526%
21	11.408	1342	1348	1355	rBV6	5283	11347	1.16%	0.100%
22	11.526	1362	1368	1374	rBV	25116	44937	4.58%	0.396%
23	11.585	1374	1378	1392	rVV	26116	53586	5.46%	0.472%
24	11.808	1410	1416	1420	rVV4	2985	6312	0.64%	0.056%
25	12.413	1511	1519	1528	rBV5	34617	85214	8.68%	0.750%
26	12.525	1534	1538	1543	rVB3	3470	5419	0.55%	0.048%
27	12.595	1545	1550	1556	rBV	2847	4686	0.48%	0.041%
28	13.118	1633	1639	1645	rBV4	1915	4767	0.49%	0.042%
29	13.183	1645	1650	1651	rVV4	3193	4755	0.48%	0.042%
30	13.283	1663	1667	1671	rVB4	2471	4224	0.43%	0.037%
31	13.330	1671	1675	1677	rBV4	2965	4194	0.43%	0.037%
32	13.418	1687	1690	1696	rVB3	4474	6609	0.67%	0.058%
33	13.547	1708	1712	1717	rBV5	4032	6758	0.69%	0.059%
34	13.618	1717	1724	1728	rVV3	8203	11924	1.21%	0.105%
35	13.671	1728	1733	1742	rVV4	6706	14947	1.52%	0.132%
36	13.765	1744	1749	1754	rVV3	6653	11501	1.17%	0.101%

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Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
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37	14.035	1787	1795	1802	rBV2	10641	17828	1.82%	0.157%
38	14.223	1821	1827	1840	rVB	362437	523866	53.38%	4.612%
39	14.446	1857	1865	1872	rBV5	11618	29061	2.96%	0.256%
40	14.528	1872	1879	1886	rVV	398817	634063	64.61%	5.582%
41	14.687	1902	1906	1909	rVB3	7001	8712	0.89%	0.077%
42	14.752	1912	1917	1922	rBV	2622	5693	0.58%	0.050%
43	14.828	1922	1930	1939	rVB	279301	459057	46.78%	4.041%
44	15.063	1964	1970	1978	rBV	24537	47374	4.83%	0.417%
45	15.551	2048	2053	2056	rBV2	8304	11398	1.16%	0.100%
46	15.586	2056	2059	2062	rVV	8833	13065	1.33%	0.115%
47	15.627	2062	2066	2071	rVB2	26319	41369	4.22%	0.364%
48	15.762	2084	2089	2091	rVV5	5166	7745	0.79%	0.068%
49	15.821	2093	2099	2106	rVB	527338	814128	82.96%	7.167%
50	15.950	2115	2121	2131	rBV2	107244	170820	17.41%	1.504%
51	16.044	2131	2137	2142	rBV5	6684	11502	1.17%	0.101%
52	16.091	2142	2145	2149	rVB	9126	11636	1.19%	0.102%
53	16.156	2152	2156	2160	rBV2	13012	15252	1.55%	0.134%
54	16.262	2170	2174	2178	rBV5	3760	5980	0.61%	0.053%
55	16.332	2181	2186	2189	rBV6	6751	11636	1.19%	0.102%
56	16.432	2197	2203	2206	rBV5	9228	16247	1.66%	0.143%
57	16.467	2206	2209	2215	rVV3	15045	28972	2.95%	0.255%
58	16.549	2221	2223	2226	rBV2	3677	5285	0.54%	0.047%
59	16.585	2226	2229	2232	rVB4	4760	6480	0.66%	0.057%
60	16.614	2232	2234	2242	rVB6	4519	9065	0.92%	0.080%
61	16.685	2242	2246	2251	rBV6	7088	12734	1.30%	0.112%
62	16.943	2287	2290	2298	rVB8	3790	7635	0.78%	0.067%
63	17.031	2301	2305	2309	rVB6	6107	8556	0.87%	0.075%
64	17.237	2334	2340	2344	rBV6	4513	8513	0.87%	0.075%
65	17.284	2344	2348	2353	rVV3	16637	26584	2.71%	0.234%
66	17.337	2354	2357	2361	rVV4	8768	13256	1.35%	0.117%
67	17.390	2361	2366	2370	rVV4	4569	10187	1.04%	0.090%
68	17.448	2372	2376	2380	rVB3	23401	27833	2.84%	0.245%
69	17.584	2392	2399	2405	rVV	326949	515515	52.53%	4.538%
70	17.678	2409	2415	2425	rVB	552957	869753	88.62%	7.657%
71	17.760	2426	2429	2434	rVB6	4265	7171	0.73%	0.063%
72	17.824	2437	2440	2442	rBV3	8047	9448	0.96%	0.083%
73	17.883	2442	2450	2454	rBV2	20222	43494	4.43%	0.383%
74	17.924	2454	2457	2458	rBV	3836	4178	0.43%	0.037%
75	18.024	2472	2474	2478	rVB5	5603	6136	0.63%	0.054%
76	18.083	2480	2484	2490	rVB3	6933	12868	1.31%	0.113%

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77	18.142	2490	2494	2499	rBV	4832	7508	0.77%	0.066%
78	18.206	2501	2505	2511	rVB2	26695	38851	3.96%	0.342%
79	18.436	2539	2544	2550	rBV2	38491	63636	6.48%	0.560%
80	18.500	2551	2555	2560	rVV2	6603	14501	1.48%	0.128%
81	18.712	2589	2591	2594	rBV4	5803	6947	0.71%	0.061%
82	18.964	2631	2634	2638	rVB5	5068	4645	0.47%	0.041%
83	19.029	2642	2645	2648	rBV4	5821	7252	0.74%	0.064%
84	19.282	2684	2688	2694	rBV3	43283	77808	7.93%	0.685%
85	19.522	2726	2729	2731	rBV3	8009	10481	1.07%	0.092%
86	19.593	2738	2741	2745	rBV3	14412	22199	2.26%	0.195%
87	19.822	2776	2780	2784	rVB6	14976	20182	2.06%	0.178%
88	19.957	2797	2803	2811	rVB	672563	981409	100.00%	8.639%
89	20.439	2879	2885	2890	rBV6	14654	27329	2.78%	0.241%
90	20.586	2908	2910	2913	rBV4	7274	5819	0.59%	0.051%
91	21.467	3055	3060	3073	rBV	64174	137899	14.05%	1.214%
92	21.879	3124	3130	3138	rVB	275428	496412	50.58%	4.370%
93	22.025	3153	3155	3161	rVB5	12562	14283	1.46%	0.126%
94	23.065	3327	3332	3337	rVB2	27282	51130	5.21%	0.450%
95	23.588	3415	3421	3428	rVB2	52467	109398	11.15%	0.963%
96	24.223	3525	3529	3539	rVB5	16995	37019	3.77%	0.326%
97	25.051	3659	3670	3684	rVB2	319042	927662	94.52%	8.166%
98	25.286	3701	3710	3724	rVB3	160269	517167	52.70%	4.553%
99	27.084	4008	4016	4035	rBV3	46645	237900	24.24%	2.094%
100	27.572	4090	4099	4122	rBV3	66909	390575	39.80%	3.438%

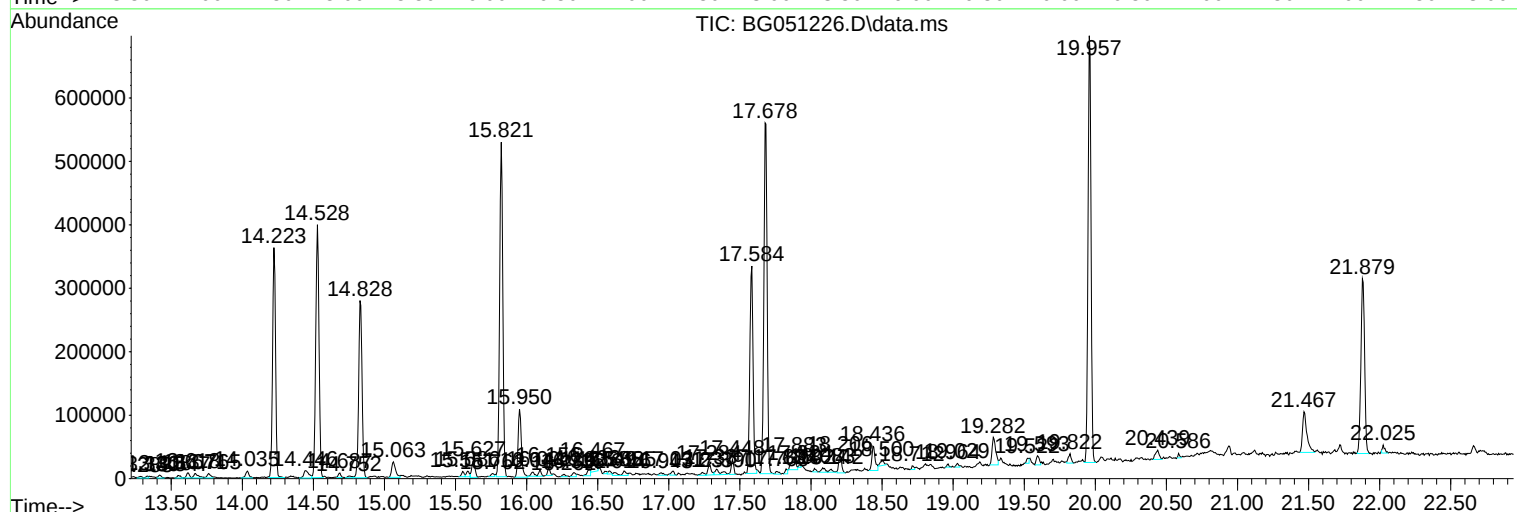
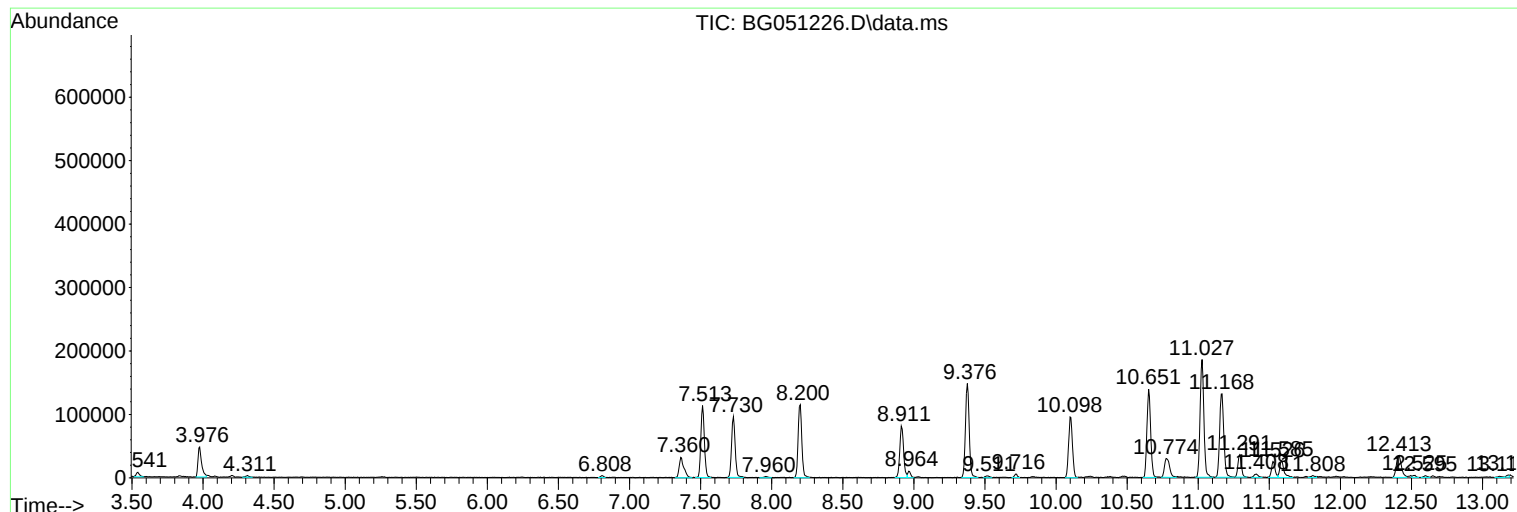
Sum of corrected areas: 11359654

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

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



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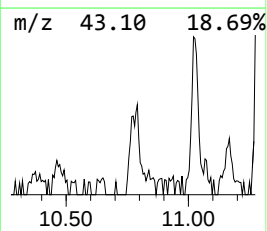
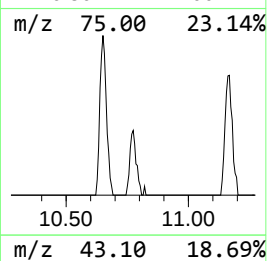
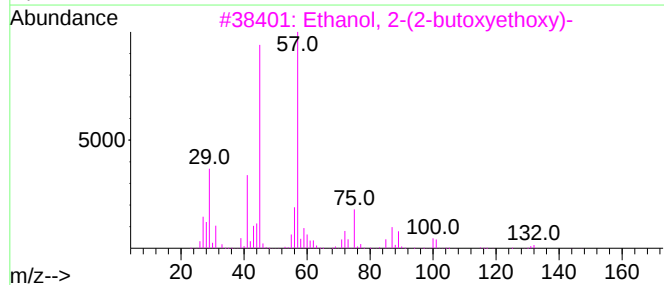
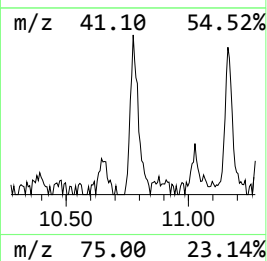
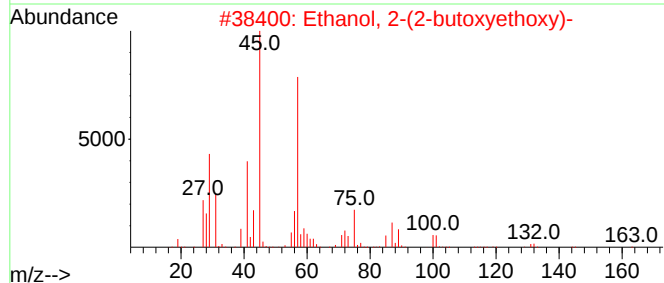
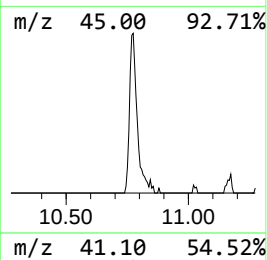
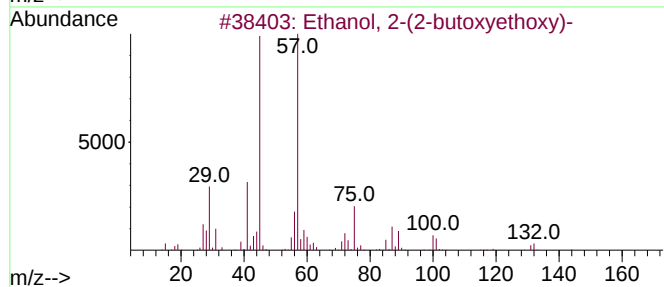
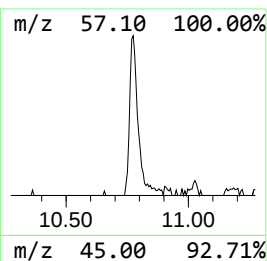
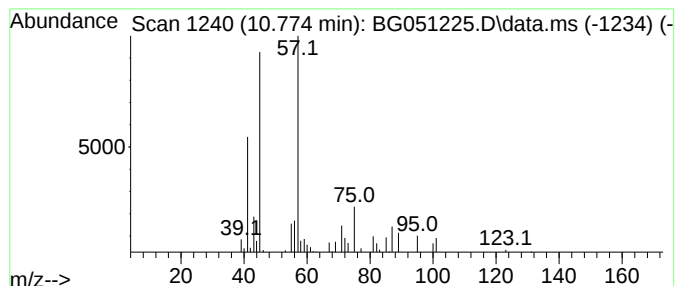
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.774	4.40 ng/ul	75344	Naphthalene-d8	11.027

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	86
2			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	78
3			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	64
4			Ethanol, 2-[2-(2-methylpropoxy)e...	162	C8H18O3	018912-80-6	53
5			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	53



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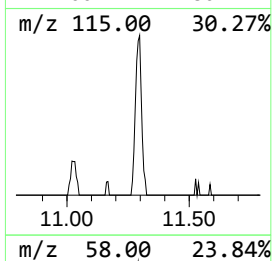
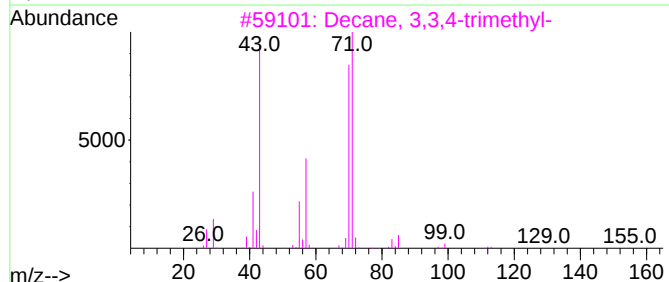
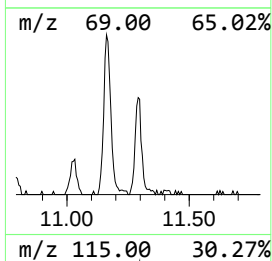
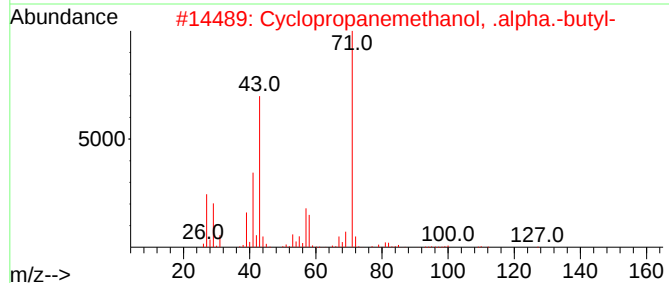
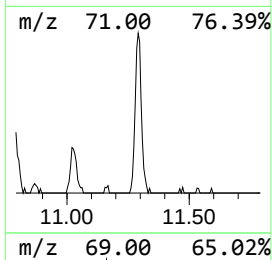
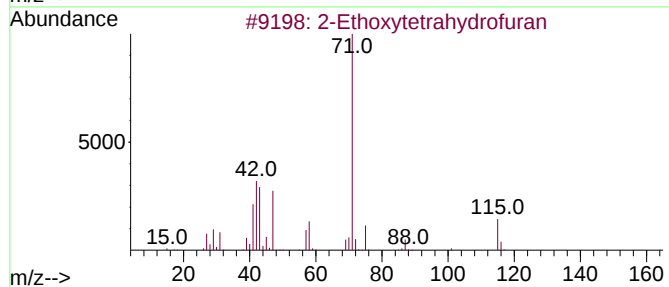
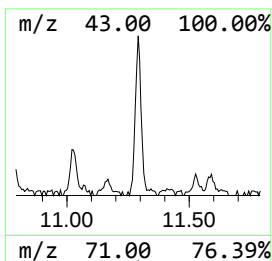
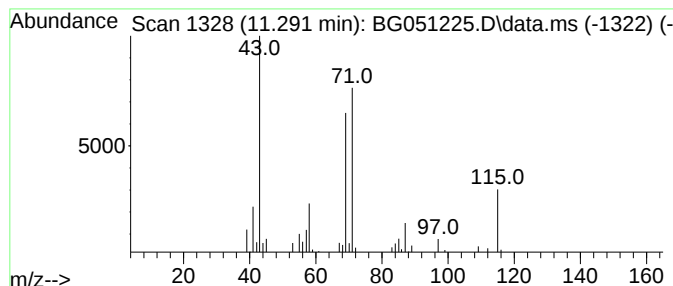
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 2-Ethoxytetrahydrofuran Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.291	3.49 ng/ul	59706	Naphthalene-d8	11.027

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Ethoxytetrahydrofuran	116	C6H12O2	013436-46-9	50
2			Cyclopropanemethanol, .alpha.-bu...	128	C8H16O	004379-16-2	47
3			Decane, 3,3,4-trimethyl-	184	C13H28	049622-18-6	43
4			Heptane, 2,5,5-trimethyl-	142	C10H22	001189-99-7	43
5			Butyric acid, 1-propylpentyl ester	200	C12H24O2	020286-46-8	43



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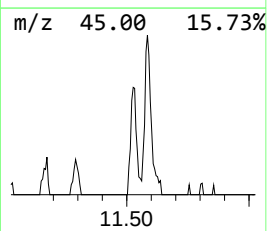
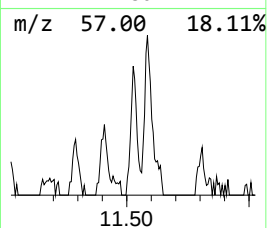
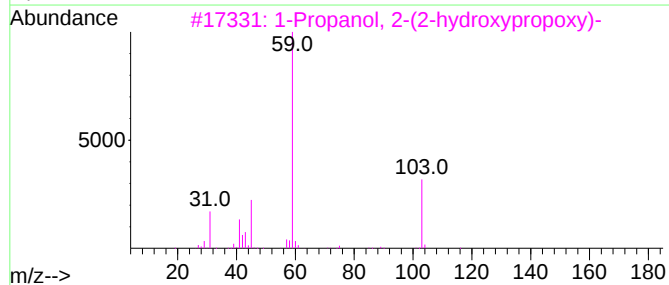
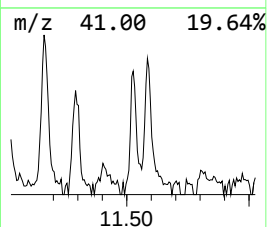
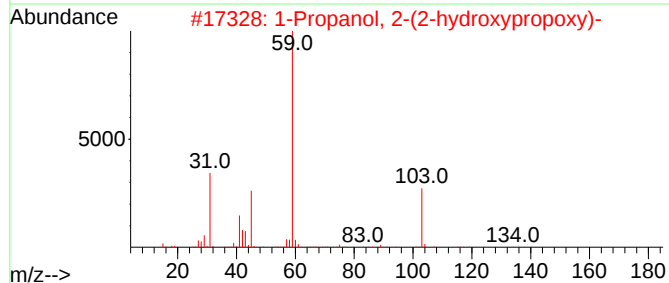
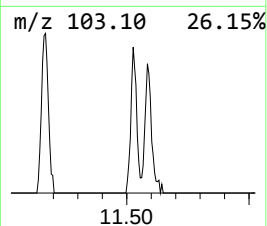
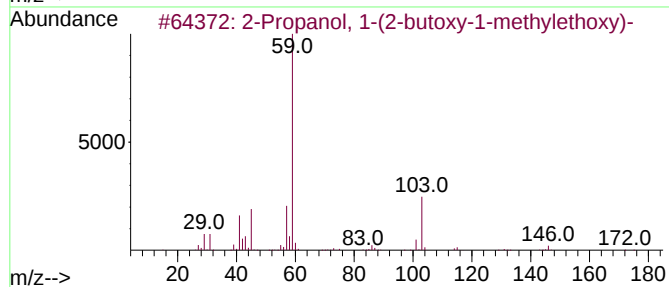
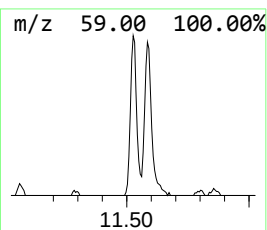
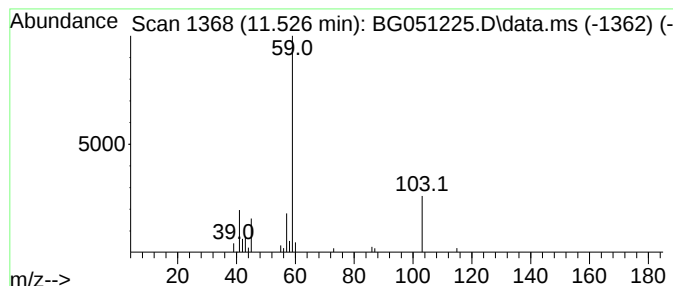
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TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 2-Propanol, 1-(2-butoxy-1-m... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.526	2.63 ng/ul	44937	Naphthalene-d8	11.027

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanol, 1-(2-butoxy-1-methyl...	190	C10H22O3	029911-28-2	83		
2	1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	78		
3	1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	78		
4	2-Propanol, 1-[1-methyl-2-(2-pro...	174	C9H18O3	055956-25-7	74		
5	1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	64		



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Data File : BG051225.D
Acq On : 25 Nov 2021 00:21
Operator : CG/JU
Sample : M4725-01
Misc :
ALS Vial : 47 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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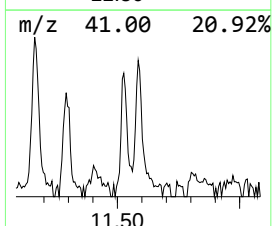
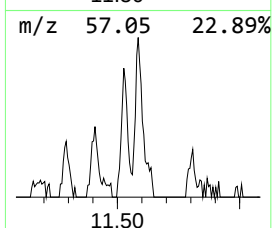
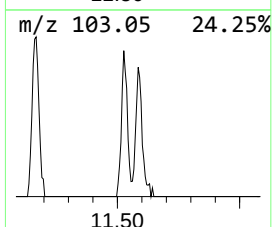
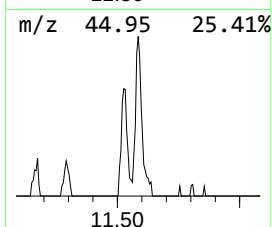
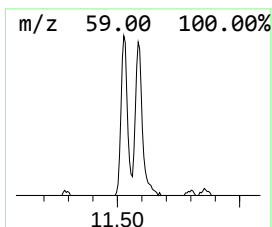
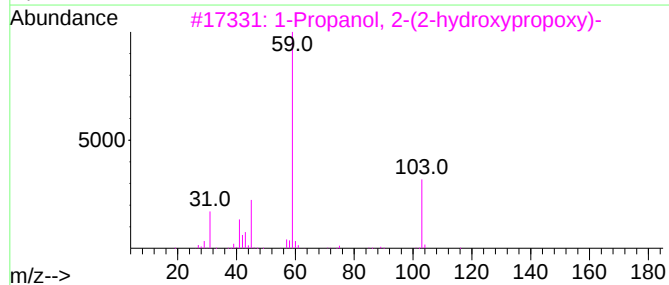
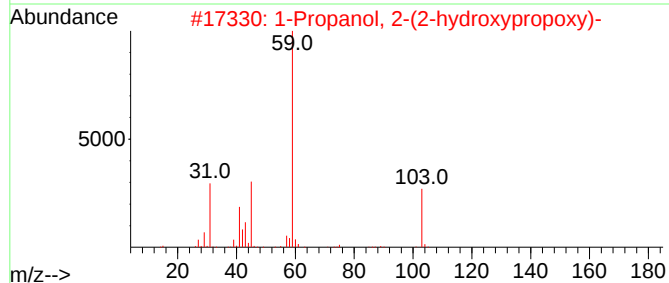
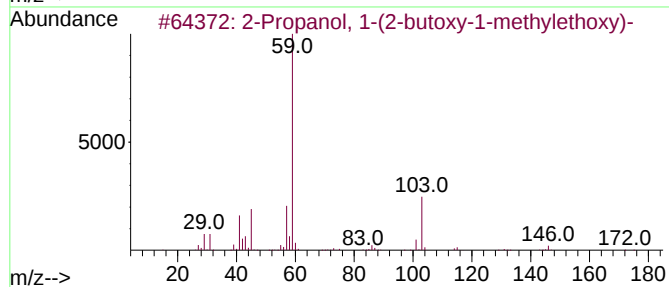
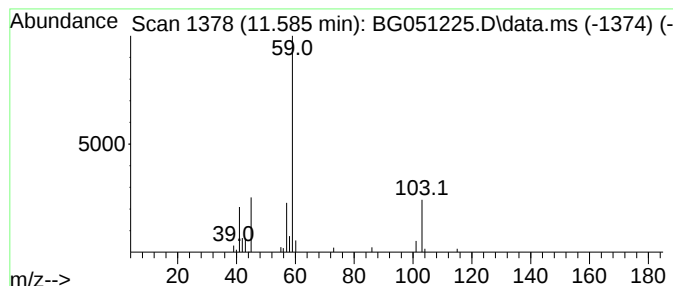
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 1-Propanol, 2-(2-hydroxypro... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.585	3.13 ng/ul	53586	Naphthalene-d8	11.027

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Propanol, 1-(2-butoxy-1-methyl...	190	C10H22O3	029911-28-2	83
2		1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	78
3		1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	78
4		1-Propanol, 2,2'-oxybis-	134	C6H14O3	000108-61-2	78
5		Dipropylene glycol (isomer 2)	134	C6H14O3	1000506-25-4	78



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG112321\
Data File : BG051225.D
Acq On : 25 Nov 2021 00:21
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Sample : M4725-01
Misc :
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Instrument :
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ClientSampleId :
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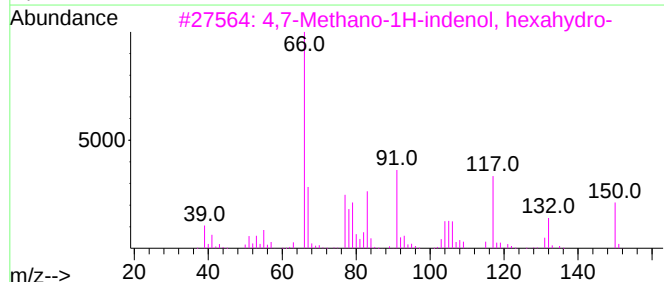
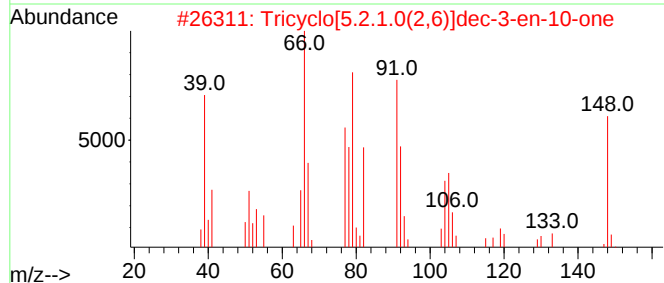
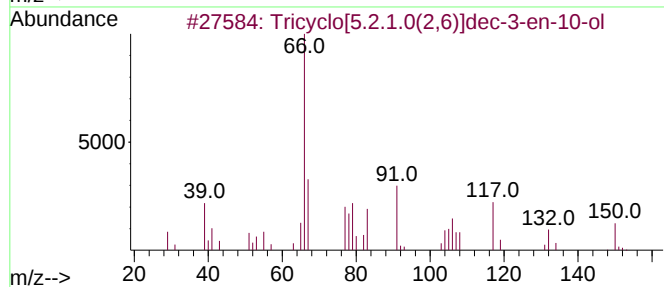
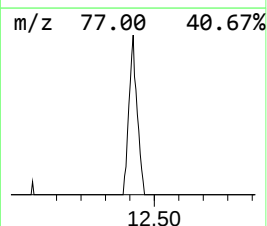
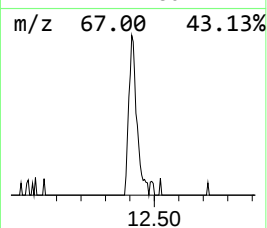
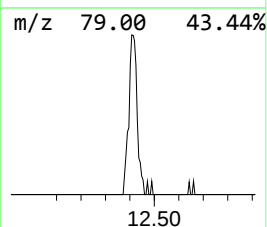
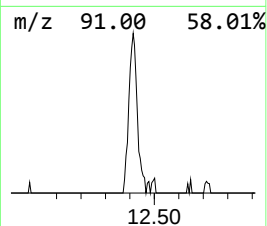
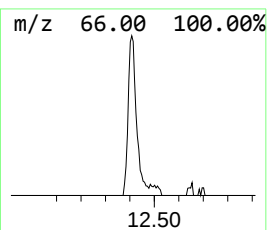
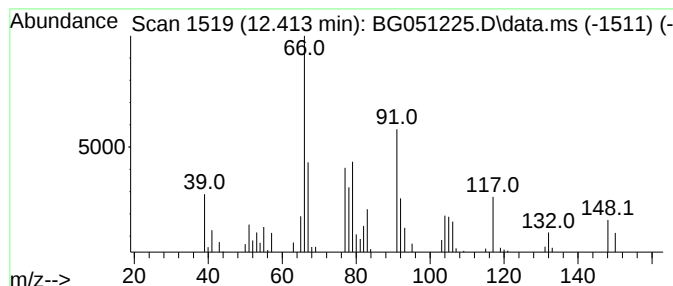
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Tricyclo[5.2.1.0(2,6)]dec-3... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.413	4.98 ng/ul	85214	Naphthalene-d8	11.027

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tricyclo[5.2.1.0(2,6)]dec-3-en-1...	150	C10H14O	039852-87-4	81
2			Tricyclo[5.2.1.0(2,6)]dec-3-en-1...	148	C10H12O	1000289-10-5	80
3			4,7-Methano-1H-indenol, hexahydro-	150	C10H14O	037275-49-3	80
4			4,7-Methano-1H-indenol, hexahydro-	150	C10H14O	037275-49-3	74
5			3a,4,5,6,7,7a-Hexahydro-4,7-meth...	134	C10H14	004488-57-7	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG112321\
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Sample : M4725-01
Misc :
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ClientSampleId :
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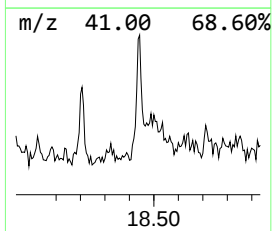
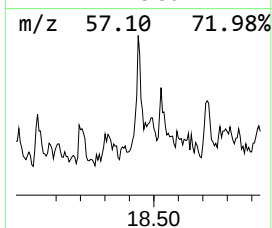
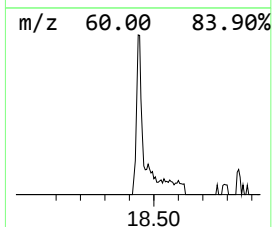
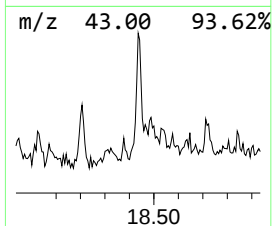
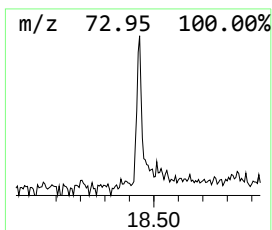
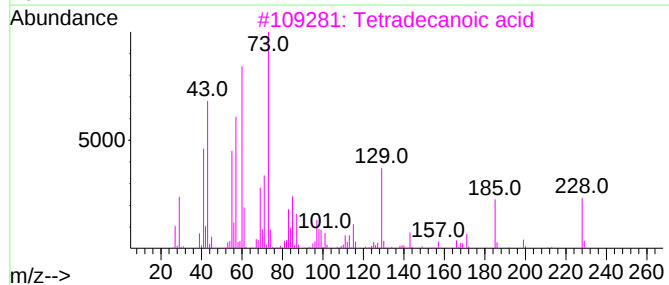
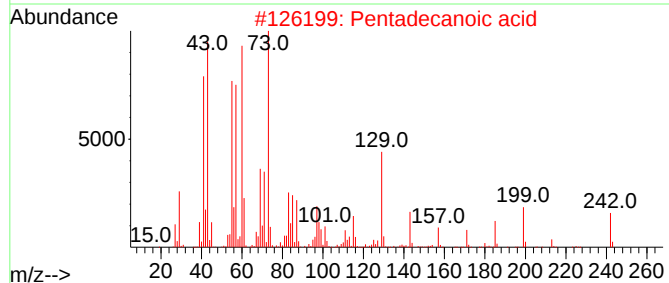
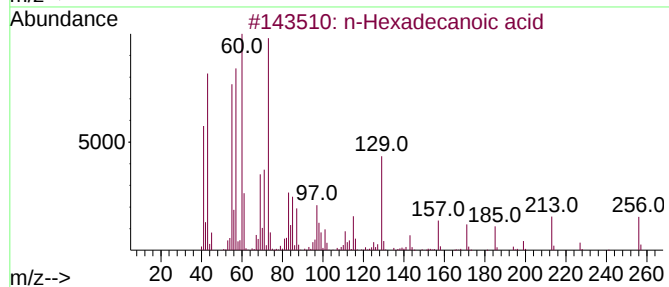
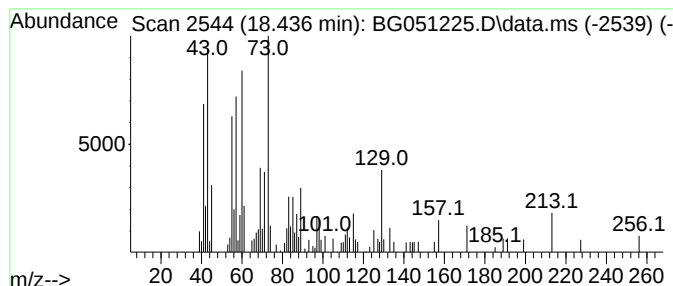
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 n-Hexadecanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.436	2.47 ng/ul	63636	Phenanthrene-d10	17.584

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	94
2			Pentadecanoic acid	242	C15H30O2	001002-84-2	81
3			Tetradecanoic acid	228	C14H28O2	000544-63-8	74
4			Tetradecanoic acid	228	C14H28O2	000544-63-8	59
5			Tridecanoic acid	214	C13H26O2	000638-53-9	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG112321\
Data File : BG051225.D
Acq On : 25 Nov 2021 00:21
Operator : CG/JU
Sample : M4725-01
Misc :
ALS Vial : 47 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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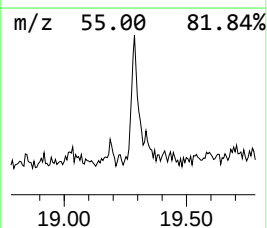
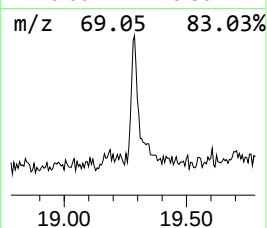
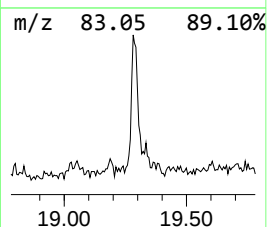
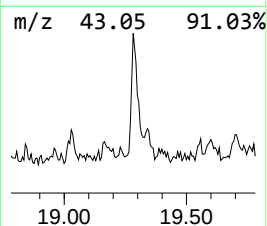
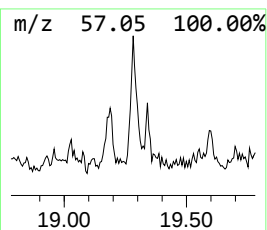
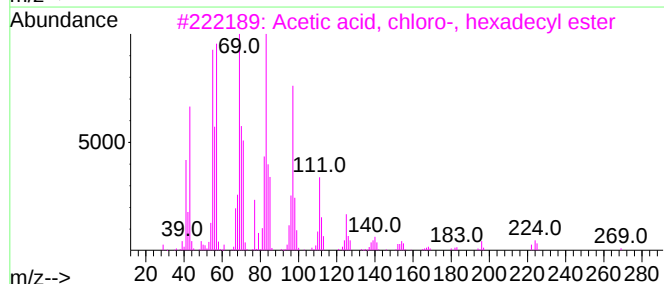
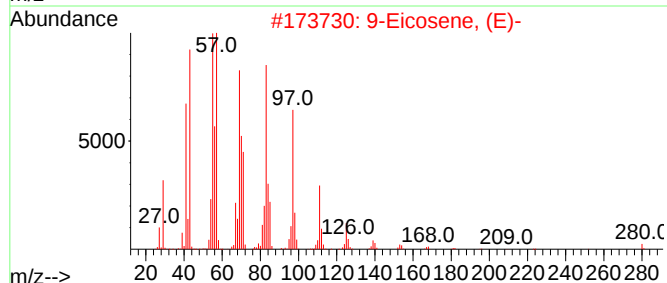
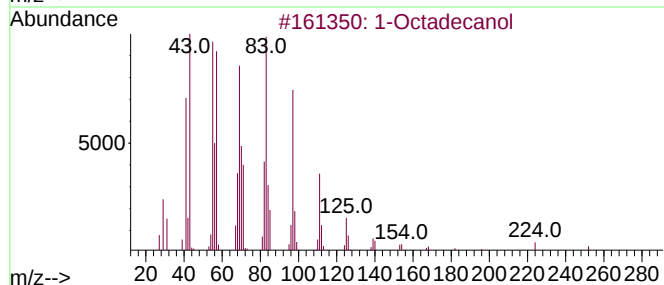
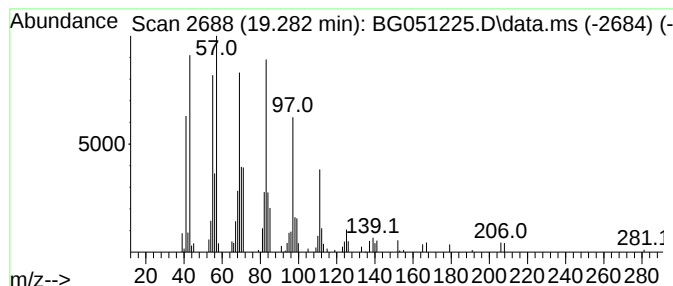
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 1-Octadecanol Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.282	3.02 ng/ul	77808	Phenanthrene-d10	17.584

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Octadecanol	270	C18H38O	000112-92-5	90
2			9-Eicosene, (E)-	280	C20H40	074685-29-3	90
3			Acetic acid, chloro-, hexadecyl ...	318	C18H35ClO2	052132-58-8	90
4			5-Octadecene, (E)-	252	C18H36	007206-21-5	87
5			Pentafluoropropionic acid, tride...	346	C16H27F5O2	959261-22-4	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG112321\
Data File : BG051225.D
Acq On : 25 Nov 2021 00:21
Operator : CG/JU
Sample : M4725-01
Misc :
ALS Vial : 47 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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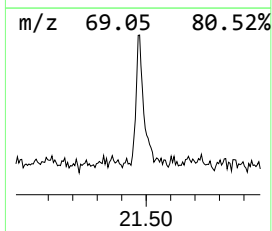
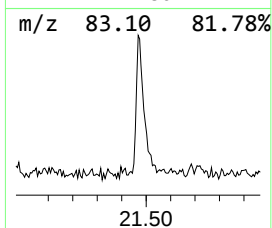
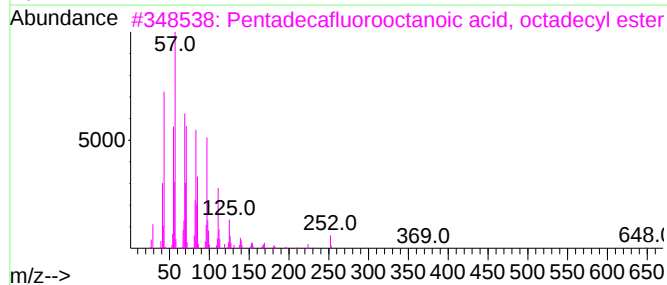
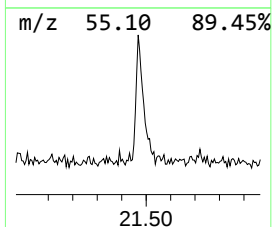
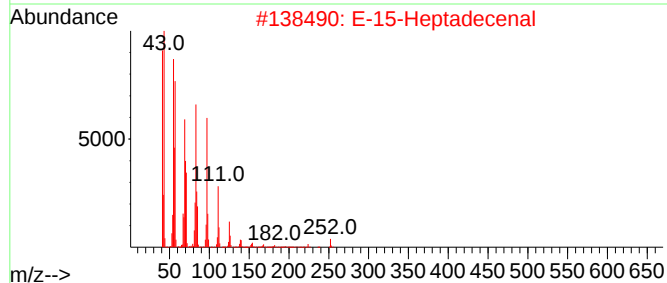
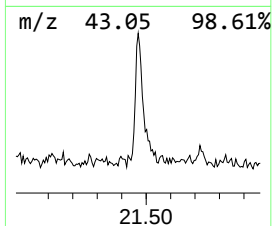
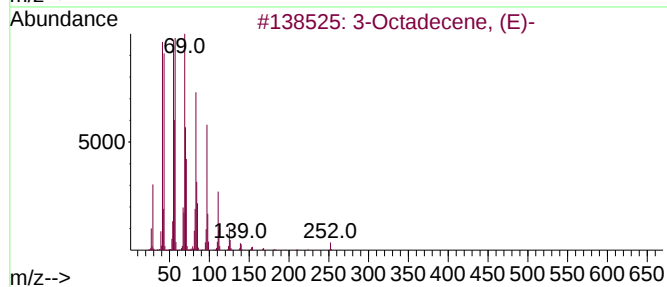
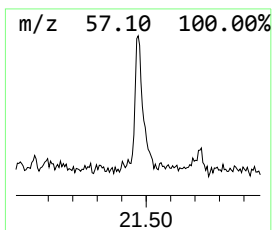
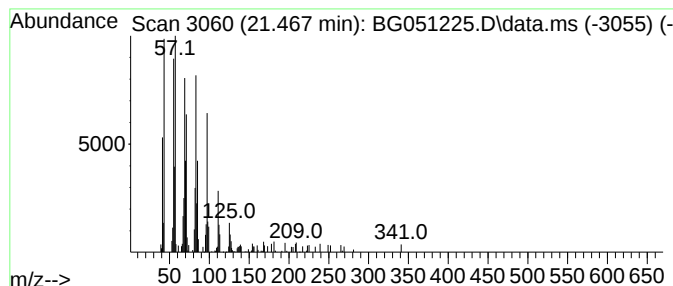
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.467	5.56 ng/ul	137899	Chrysene-d12	21.878

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Octadecene, (E)-	252	C18H36	007206-19-1	96
2			E-15-Heptadecenal	252	C17H32O	1000130-97-9	92
3			Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	91
4			Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	90
5			Pentafluoropropionic acid, penta...	374	C18H31F5O2	959092-08-1	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG112321\
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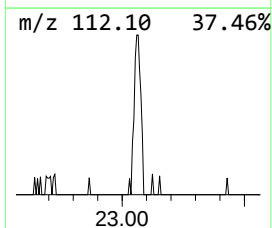
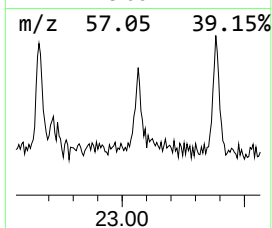
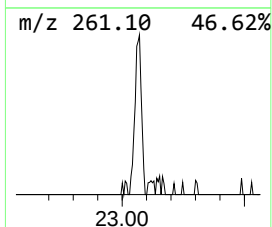
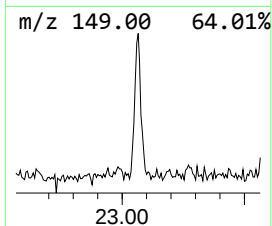
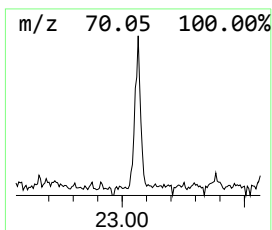
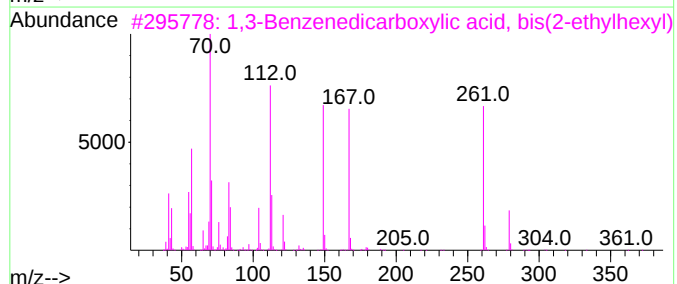
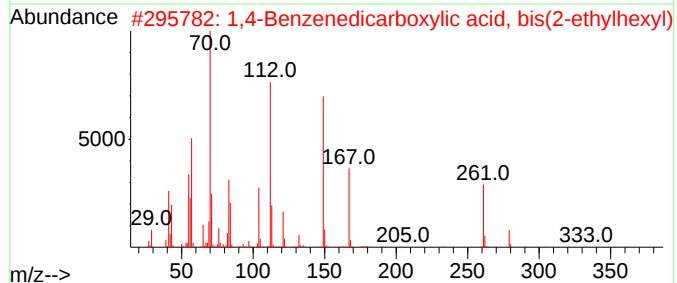
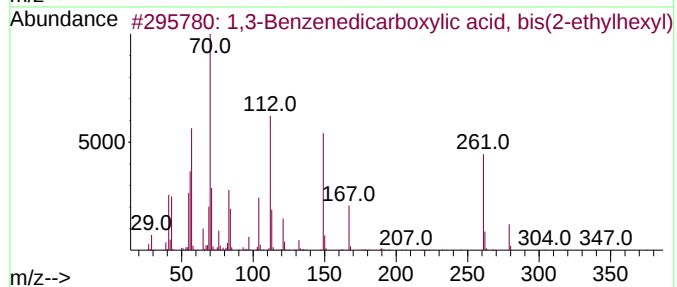
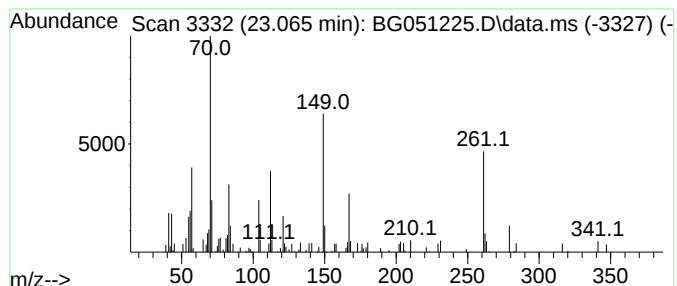
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 1,3-Benzenedicarboxylic aci... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
23.065	2.06 ng/ul	51130	Chrysene-d12	21.878	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Benzenedicarboxylic acid, bi...	390	C24H38O4	000137-89-3	80
2	1,4-Benzenedicarboxylic acid, bi...	390	C24H38O4	006422-86-2	59
3	1,3-Benzenedicarboxylic acid, bi...	390	C24H38O4	000137-89-3	43
4	4-Nitrobenzoic acid, cyclobutyl ...	221	C11H11NO4	070335-00-1	35
5	Cyclohexanamine, N-but-2-enylid...	167	C10H17NO	068048-01-1	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG112321\
Data File : BG051225.D
Acq On : 25 Nov 2021 00:21
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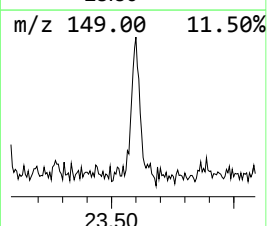
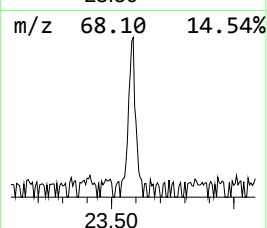
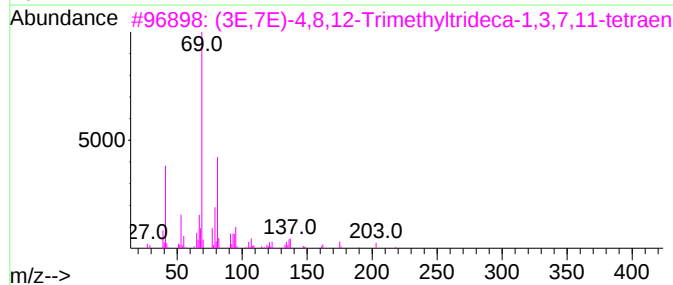
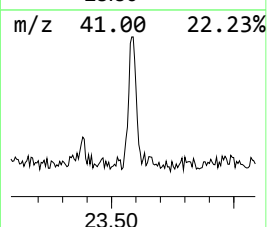
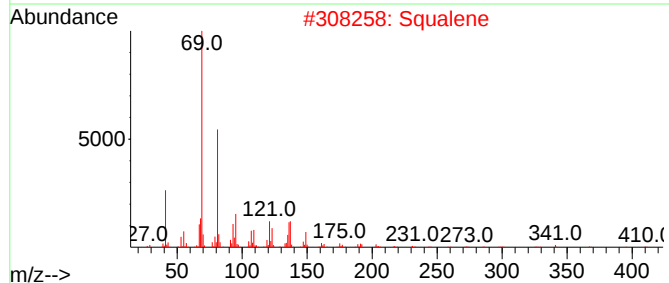
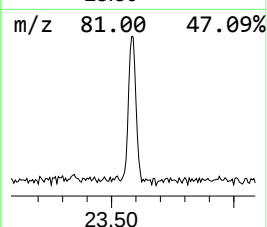
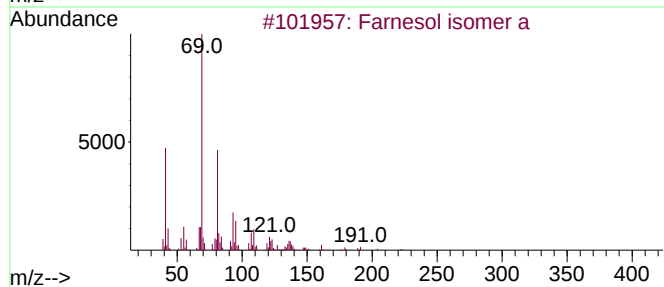
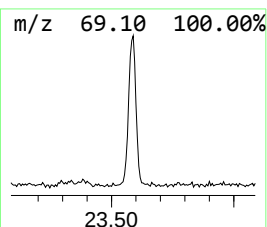
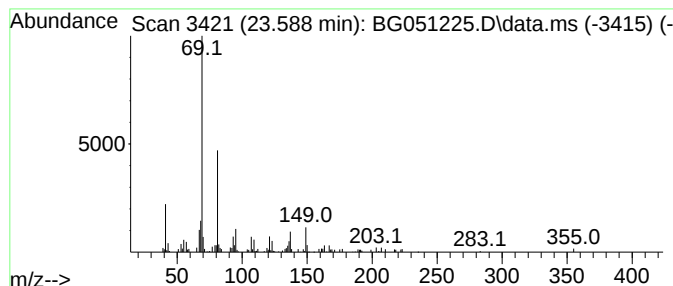
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Farnesol isomer a Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.588	4.23 ng/ul	109398	Perylene-d12	25.286

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Farnesol isomer a	222	C15H26O	1000108-92-4	81
2			Squalene	410	C30H50	000111-02-4	72
3			(3E,7E)-4,8,12-Trimethyltrideca...	218	C16H26	062235-06-7	68
4			Hexadeca-2,6,10,14-tetraen-1-ol,...	290	C20H34O	007614-21-3	64
5			3,7,11-Trimethyl-2,6,10-dodecatr...	217	C15H23N	029789-67-1	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG112321\
Data File : BG051225.D
Acq On : 25 Nov 2021 00:21
Operator : CG/JU
Sample : M4725-01
Misc :
ALS Vial : 47 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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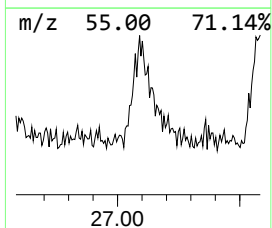
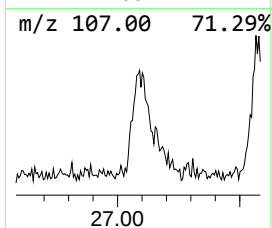
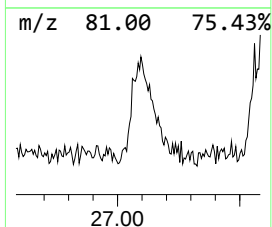
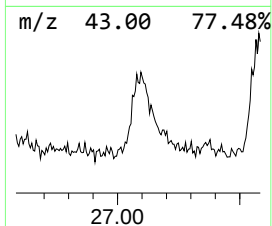
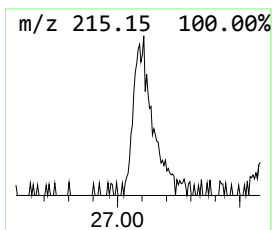
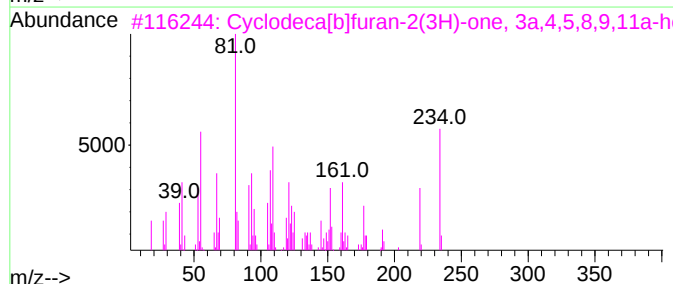
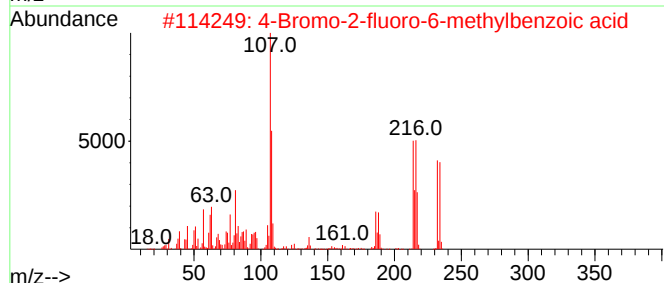
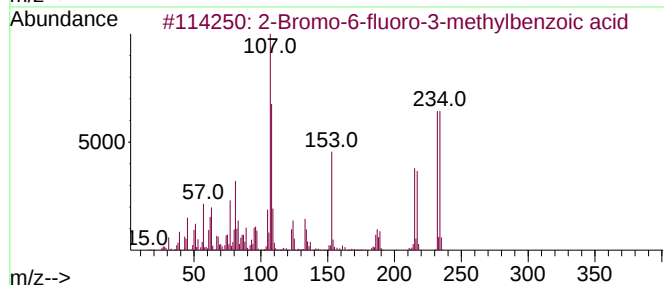
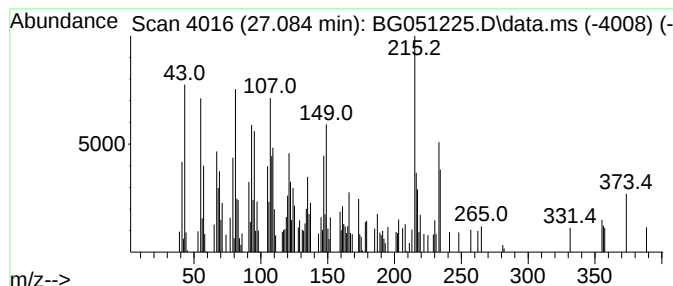
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 unknown-01 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
27.084	9.20 ng/ul	237900	Perylene-d12	25.286

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Bromo-6-fluoro-3-methylbenzoic...	232	C8H6BrF02	1359857-60-5	27		
2	4-Bromo-2-fluoro-6-methylbenzoic...	232	C8H6BrF02	1242157-23-8	27		
3	Cyclodeca[b]furan-2(3H)-one, 3a,...	234	C15H22O2	002225-79-8	25		
4	2-Norbornanamine, N-ethyl-3-phenyl-	215	C15H21N	001209-98-9	15		
5	Phthalic acid, 4-fluorobenzyl he...	358	C21H23F04	1000377-74-5	14		



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG112321\
Data File : BG051225.D
Acq On : 25 Nov 2021 00:21
Operator : CG/JU
Sample : M4725-01
Misc :
ALS Vial : 47 Sample Multiplier: 1

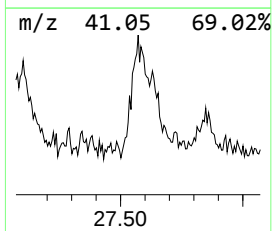
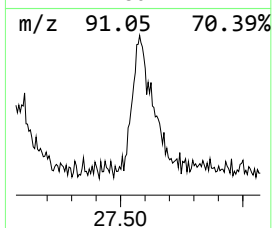
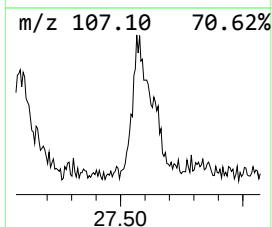
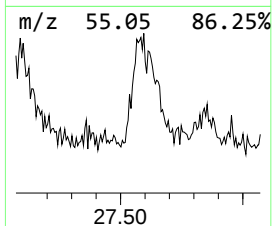
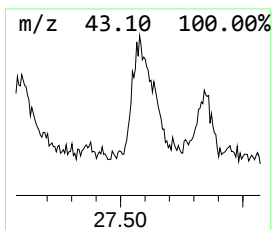
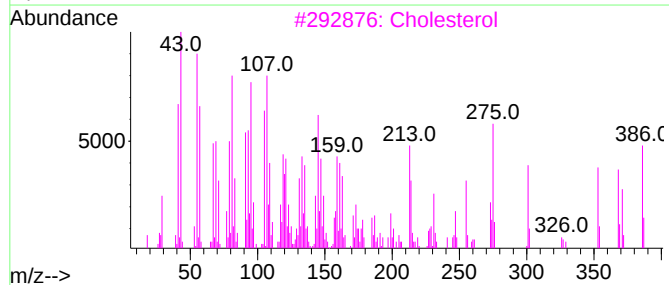
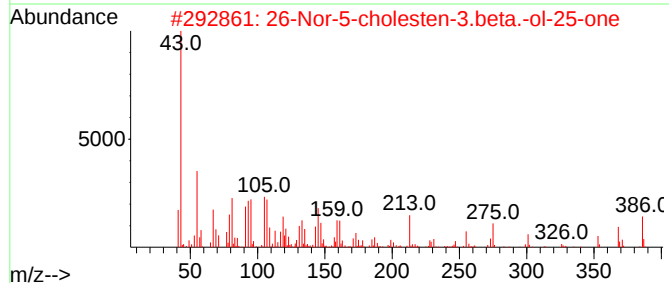
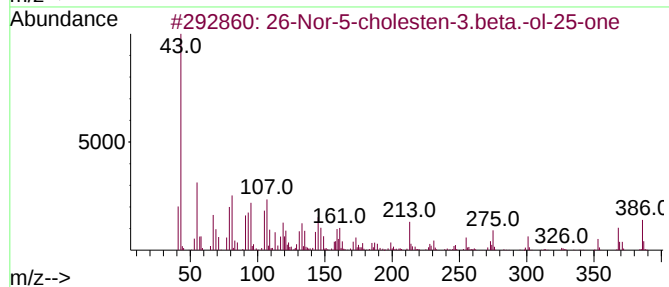
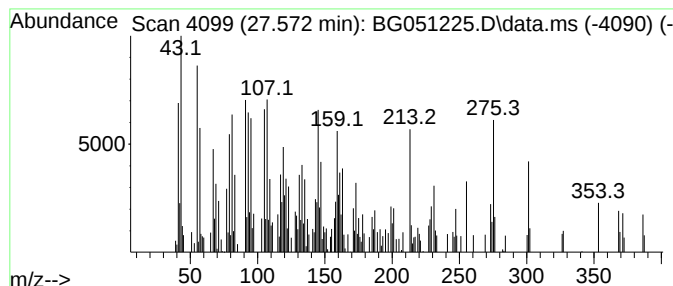
Instrument :
BNA_G
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

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

Peak Number 12 26-Nor-5-cholesten-3.beta.-... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD		R.T.	
27.572	15.10 ng/ul	390575	Perylene-d12		25.286	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	26-Nor-5-cholesten-3.beta.-ol-25...		386	C26H42O2	007494-34-0	97
2	26-Nor-5-cholesten-3.beta.-ol-25...		386	C26H42O2	007494-34-0	89
3	Cholesterol		386	C27H46O	000057-88-5	87
4	Cholesterol		386	C27H46O	000057-88-5	86
5	Propanamide, 3-phenyl-N-nonyl-		275	C18H29NO	1000407-15-6	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG112321\
Data File : BG051225.D
Acq On : 25 Nov 2021 00:21
Operator : CG/JU
Sample : M4725-01
Misc :
ALS Vial : 47 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
F4L00

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.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
Ethanol, 2-(2-b...	10.774	4.4	ng/ul	75344	2	11.027	342255	20.0
2-Ethoxytetrahy...	11.291	3.5	ng/ul	59706	2	11.027	342255	20.0
2-Propanol, 1-(...	11.526	2.6	ng/ul	44937	2	11.027	342255	20.0
1-Propanol, 2-(...	11.585	3.1	ng/ul	53586	2	11.027	342255	20.0
Tricyclo[5.2.1...	12.413	5.0	ng/ul	85214	2	11.027	342255	20.0
n-Hexadecanoic ...	18.436	2.5	ng/ul	63636	4	17.584	515515	20.0
1-Octadecanol	19.282	3.0	ng/ul	77808	4	17.584	515515	20.0
(DEL) Alkane: S...	21.467	5.6	ng/ul	137899	5	21.878	496412	20.0
1,3-Benzenedica...	23.065	2.1	ng/ul	51130	5	21.878	496412	20.0
Farnesol isomer a	23.588	4.2	ng/ul	109398	6	25.286	517167	20.0
unknown-01	27.084	9.2	ng/ul	237900	6	25.286	517167	20.0
26-Nor-5-choles...	27.572	15.1	ng/ul	390575	6	25.286	517167	20.0