

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072621\
 Data File : BG049522.D
 Acq On : 28 Jul 2021 00:51
 Operator : CG/JU
 Sample : M3091-01
 Misc :
 ALS Vial : 52 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0036

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-BG072421.M
 Title : SVOA CALIBRATION

Signal : TIC: BG049522.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.078	3	7	17	rBV2	39004	75427	12.72%	1.008%
2	3.161	17	21	28	rVB	23418	35022	5.91%	0.468%
3	3.425	63	66	73	rVB4	3523	5633	0.95%	0.075%
4	3.860	136	140	151	rBV2	26818	57556	9.71%	0.769%
5	4.970	326	329	332	rBV2	1163	1826	0.31%	0.024%
6	7.202	702	709	720	rVB2	90411	162128	27.34%	2.166%
7	7.367	731	737	743	rBV	55551	91118	15.36%	1.217%
8	7.578	766	773	782	rVB	84258	149761	25.25%	2.001%
9	8.048	846	853	860	rVV	79191	132441	22.33%	1.769%
10	8.753	967	973	981	rVB2	115524	195522	32.97%	2.612%
11	9.211	1041	1051	1062	rVB	91423	168422	28.40%	2.250%
12	9.382	1076	1080	1085	rBV2	1577	3128	0.53%	0.042%
13	9.940	1167	1175	1182	rBV2	86479	152634	25.74%	2.039%
14	10.486	1257	1268	1277	rBV	120822	221885	37.41%	2.964%
15	10.627	1283	1292	1306	rBV	66412	123996	20.91%	1.657%
16	10.868	1326	1333	1340	rVB	110965	194762	32.84%	2.602%
17	10.997	1349	1355	1364	rBV	105142	201481	33.97%	2.692%
18	12.448	1600	1602	1609	rVB3	2106	2949	0.50%	0.039%
19	13.294	1741	1746	1750	rBV	2448	3154	0.53%	0.042%
20	13.535	1784	1787	1788	rVV	3128	2744	0.46%	0.037%
21	13.576	1788	1794	1801	rVV2	38698	63511	10.71%	0.849%
22	13.887	1844	1847	1853	rVB2	1299	2089	0.35%	0.028%
23	14.093	1876	1882	1890	rBV	250933	372050	62.74%	4.971%
24	14.328	1918	1922	1925	rBV	2613	3204	0.54%	0.043%
25	14.393	1926	1933	1941	rVB	253061	409849	69.11%	5.476%
26	14.698	1978	1985	1993	rVB	190620	303346	51.15%	4.053%
27	14.892	2012	2018	2027	rBV2	103858	182601	30.79%	2.440%
28	15.097	2047	2053	2054	rBV3	2194	3472	0.59%	0.046%
29	15.473	2113	2117	2124	rBB	1489	3560	0.60%	0.048%
30	15.532	2124	2127	2131	rBV3	2576	3660	0.62%	0.049%
31	15.691	2146	2154	2161	rVB	337711	527368	88.93%	7.046%
32	15.808	2168	2174	2180	rBV	106051	166201	28.02%	2.220%
33	15.926	2191	2194	2199	rVB3	4915	6657	1.12%	0.089%
34	16.131	2226	2229	2231	rVV2	2432	2737	0.46%	0.037%
35	16.155	2231	2233	2237	rVB4	1931	1867	0.31%	0.025%
36	16.472	2280	2287	2292	rBV5	7592	13695	2.31%	0.183%

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Smoothing : OFF

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Peak Location: TOP

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

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Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG072421.M
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37	16.572	2299	2304	2310	rVB2	3270	6130	1.03%	0.082%
38	16.672	2318	2321	2325	rVB4	3045	4545	0.77%	0.061%
39	16.825	2339	2347	2352	rBV2	5001	9612	1.62%	0.128%
40	16.895	2354	2359	2364	rBV3	3698	6907	1.16%	0.092%
41	17.136	2395	2400	2401	rBV	2567	2721	0.46%	0.036%
42	17.183	2404	2408	2411	rVV2	4949	8193	1.38%	0.109%
43	17.236	2415	2417	2419	rVV	2392	2470	0.42%	0.033%
44	17.353	2432	2437	2438	rBV	2029	2973	0.50%	0.040%
45	17.447	2446	2453	2459	rBV	248759	371943	62.72%	4.969%
46	17.547	2459	2470	2480	rVB	361587	570558	96.21%	7.623%
47	17.682	2489	2493	2497	rBV2	2857	3263	0.55%	0.044%
48	17.735	2497	2502	2503	rBV3	1344	1973	0.33%	0.026%
49	17.817	2511	2516	2519	rBV4	3494	6400	1.08%	0.086%
50	17.976	2540	2543	2545	rVB2	2353	1837	0.31%	0.025%
51	18.041	2551	2554	2556	rBV4	1916	2368	0.40%	0.032%
52	18.099	2560	2564	2570	rVB2	22455	30342	5.12%	0.405%
53	18.229	2584	2586	2591	rVB2	4155	5441	0.92%	0.073%
54	18.328	2598	2603	2608	rVV4	16274	23835	4.02%	0.318%
55	18.599	2647	2649	2650	rBV2	2516	2065	0.35%	0.028%
56	18.628	2650	2654	2659	rVV8	6896	11433	1.93%	0.153%
57	18.681	2659	2663	2665	rVV4	1456	2163	0.36%	0.029%
58	18.751	2673	2675	2677	rBV3	2926	1832	0.31%	0.024%
59	18.875	2694	2696	2697	rBV	3081	2451	0.41%	0.033%
60	18.928	2700	2705	2708	rBV4	5951	10494	1.77%	0.140%
61	18.957	2708	2710	2713	rVV3	1771	2152	0.36%	0.029%
62	19.022	2717	2721	2722	rBV3	2145	2923	0.49%	0.039%
63	19.086	2730	2732	2737	rVB3	8021	10206	1.72%	0.136%
64	19.127	2737	2739	2742	rBV4	1969	2549	0.43%	0.034%
65	19.180	2744	2748	2754	rBV4	11760	18953	3.20%	0.253%
66	19.239	2754	2758	2760	rVV3	14969	21132	3.56%	0.282%
67	19.268	2760	2763	2769	rVB	88509	104893	17.69%	1.401%
68	19.398	2781	2785	2790	rVB2	31856	39389	6.64%	0.526%
69	19.468	2795	2797	2799	rBV3	3440	2623	0.44%	0.035%
70	19.603	2816	2820	2824	rBV6	4286	7718	1.30%	0.103%
71	19.721	2836	2840	2842	rBV3	2922	4004	0.68%	0.053%
72	19.832	2853	2859	2866	rVB	410192	593048	100.00%	7.923%
73	20.126	2905	2909	2910	rBV4	2611	2804	0.47%	0.037%
74	20.320	2940	2942	2945	rBV3	4834	5456	0.92%	0.073%
75	20.520	2974	2976	2981	rVB5	4707	5562	0.94%	0.074%
76	20.590	2986	2988	2990	rBV2	2332	2245	0.38%	0.030%

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Integrator: RTE
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 Sampling : 1 Min Area: 0.1 % of largest Peak
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If leading or trailing edge < 100 prefer < Baseline drop else tangent >
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77	20.684	3000	3004	3008	rVB4	9456	12738	2.15%	0.170%
78	20.725	3008	3011	3015	rBV5	4703	5860	0.99%	0.078%
79	21.172	3085	3087	3088	rBV2	3916	2091	0.35%	0.028%
80	21.354	3114	3118	3124	rBV3	30819	52074	8.78%	0.696%
81	21.671	3170	3172	3173	rBV2	3585	2611	0.44%	0.035%
82	21.730	3175	3182	3190	rVB	223632	377255	63.61%	5.040%
83	22.352	3287	3288	3293	rBV5	4416	5135	0.87%	0.069%
84	22.957	3389	3391	3394	rVB4	4031	3232	0.54%	0.043%
85	23.210	3430	3434	3436	rBV5	9908	15805	2.67%	0.211%
86	24.062	3574	3579	3585	rVB6	10814	21694	3.66%	0.290%
87	24.309	3619	3621	3625	rBV5	3906	3998	0.67%	0.053%
88	24.408	3636	3638	3639	rBV2	3757	2209	0.37%	0.030%
89	24.778	3688	3701	3711	rBV	198419	574878	96.94%	7.680%
90	25.002	3730	3739	3750	rVB2	131196	392310	66.15%	5.241%
91	25.413	3807	3809	3812	rBV4	3788	5017	0.85%	0.067%
92	25.554	3830	3833	3838	rBV7	5216	8748	1.48%	0.117%
93	25.924	3894	3896	3897	rBV2	3385	2407	0.41%	0.032%
94	26.094	3923	3925	3926	rVB2	4706	2396	0.40%	0.032%
95	26.177	3933	3939	3940	rBV6	10216	16316	2.75%	0.218%
96	27.522	4166	4168	4169	rBV	3392	2300	0.39%	0.031%
97	27.968	4242	4244	4246	rBV3	3088	2433	0.41%	0.033%
98	30.999	4758	4760	4762	rBV3	4022	4316	0.73%	0.058%
99	31.516	4846	4848	4849	rBV2	3364	2169	0.37%	0.029%
100	31.975	4924	4926	4928	rBV3	3084	1969	0.33%	0.026%

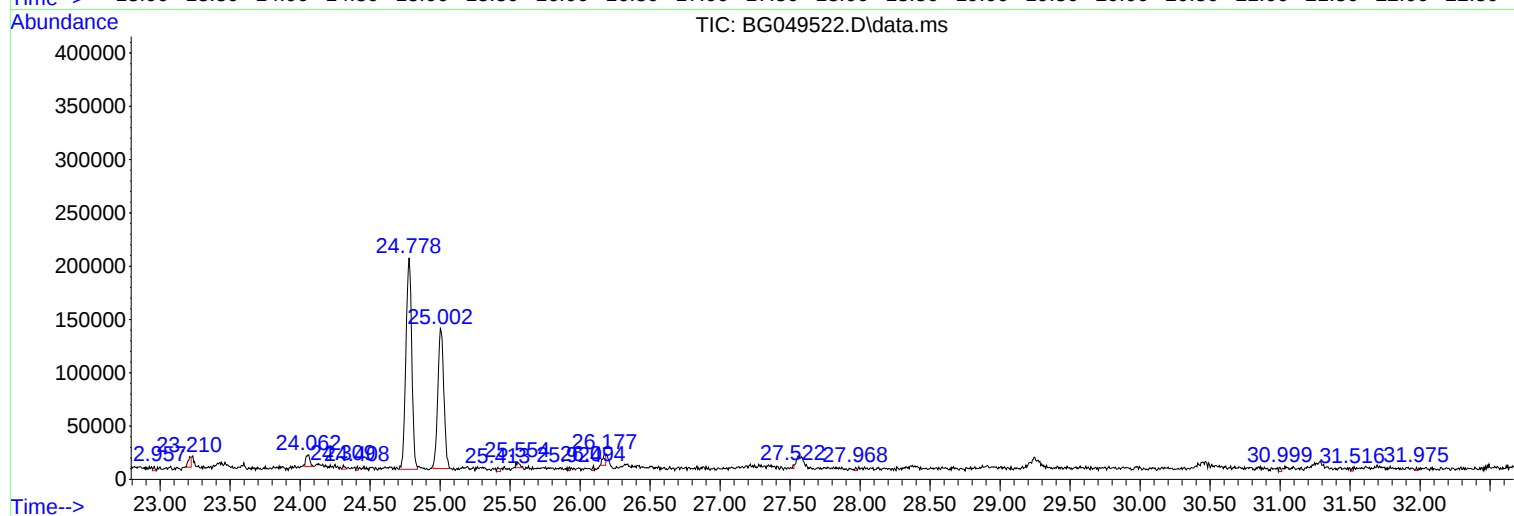
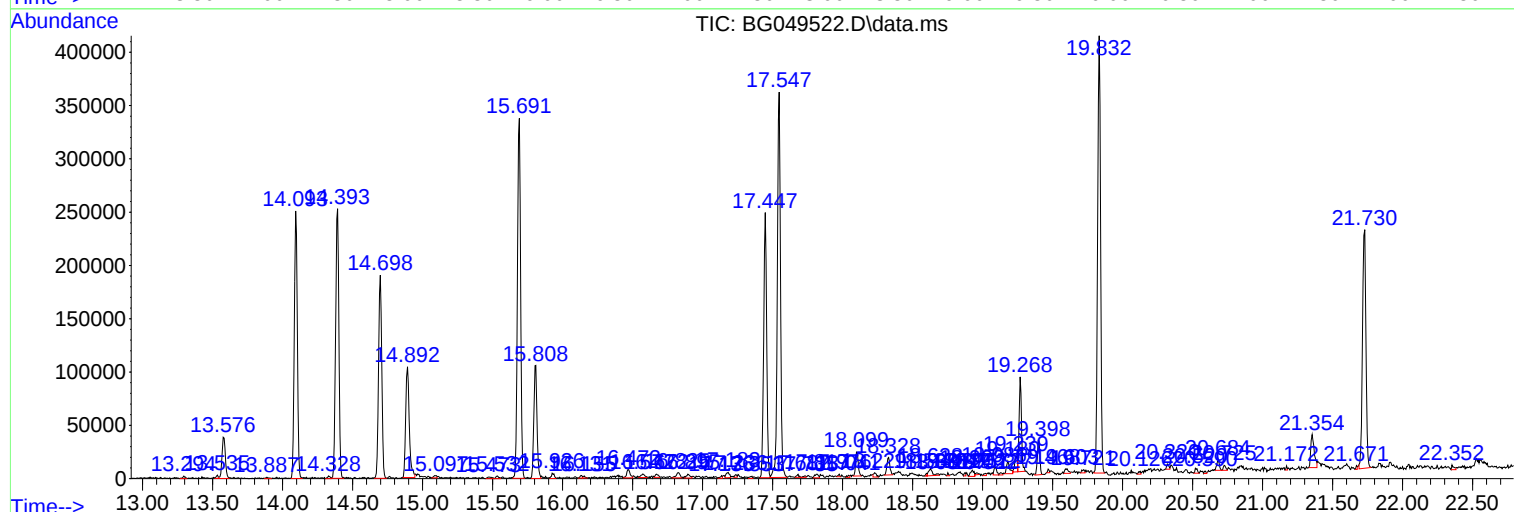
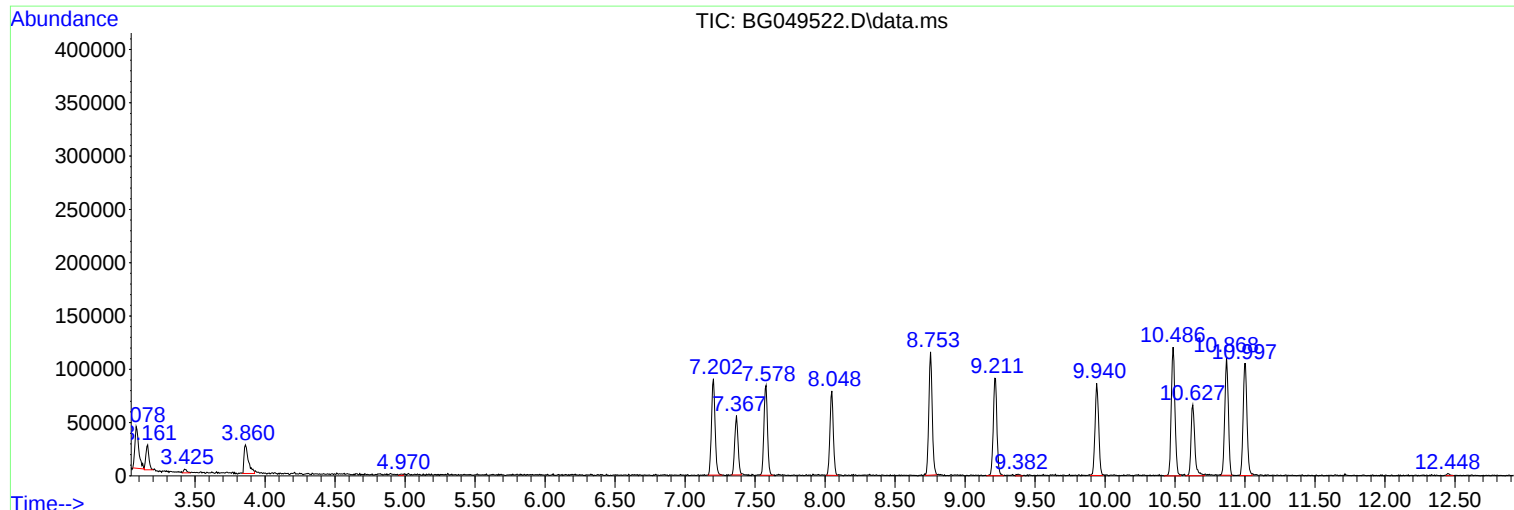
Sum of corrected areas: 7485023

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Instrument :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG072421.M
Quant Title : SVOA CALIBRATION

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



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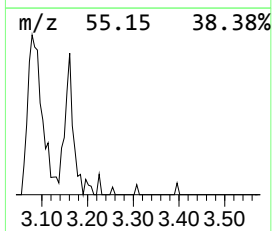
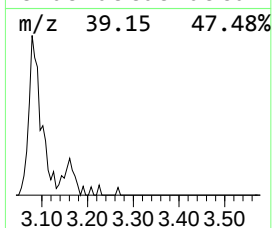
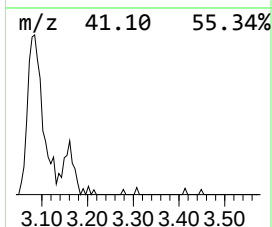
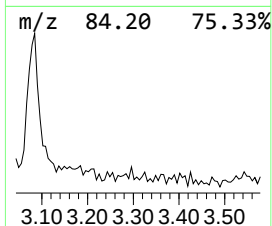
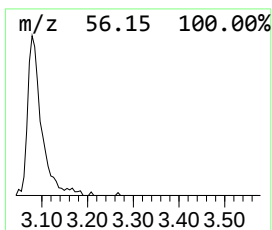
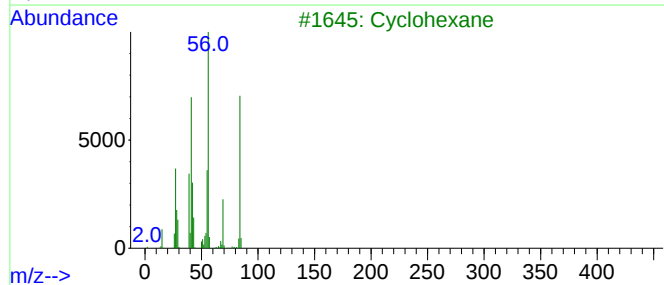
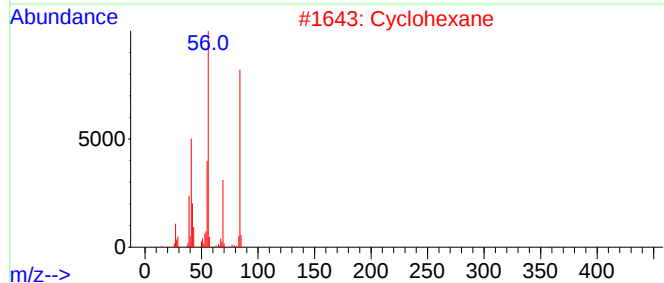
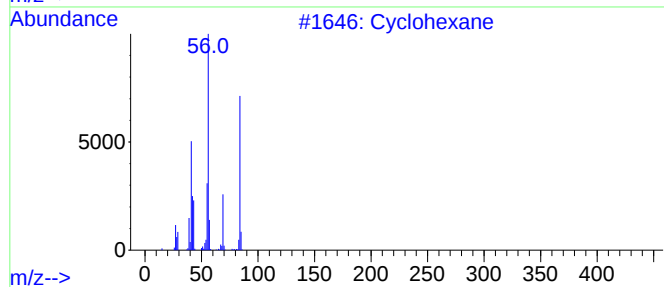
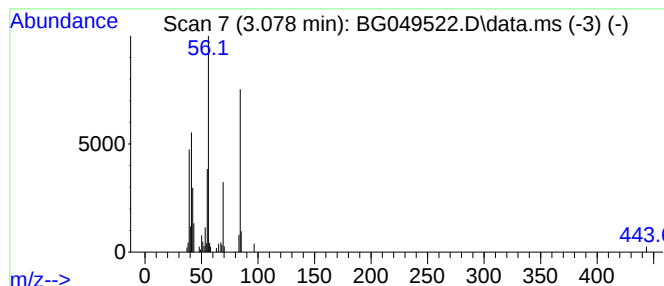
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG072421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 (DEL) Alkane: Cyclic3.078 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.078	11.39 ng/ul	75427	1,4-Dichlorobenzene-d4	8.048

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane	84	C6H12	000110-82-7	76
2			Cyclohexane	84	C6H12	000110-82-7	76
3			Cyclohexane	84	C6H12	000110-82-7	74
4			1-Pentene, 2-methyl-	84	C6H12	000763-29-1	64
5			Cyclohexane	84	C6H12	000110-82-7	64



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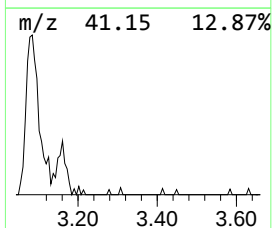
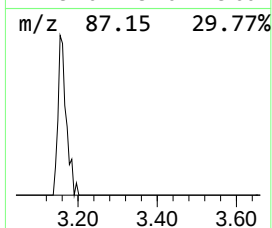
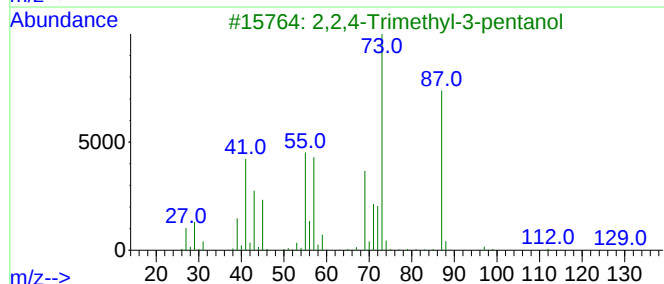
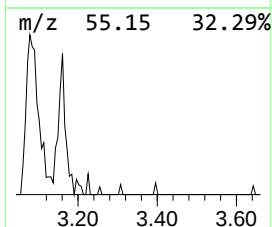
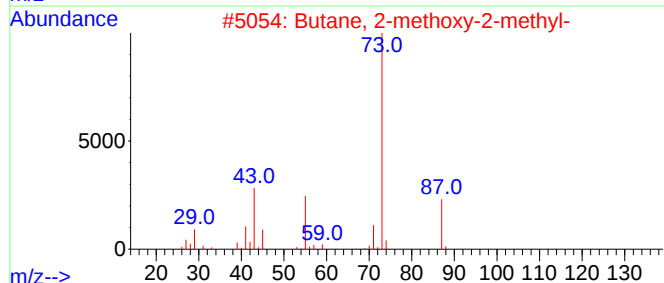
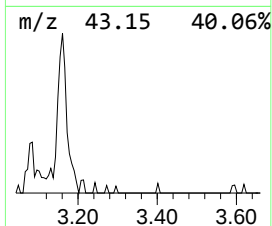
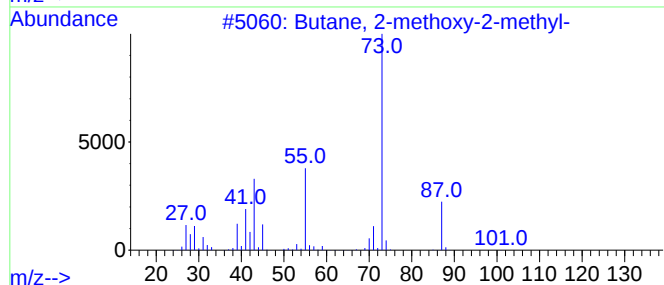
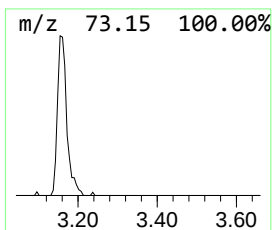
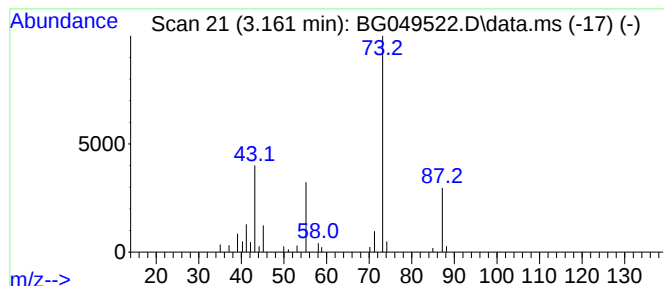
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Quant Title : SVOA CALIBRATION

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

Peak Number 2 Butane, 2-methoxy-2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.161	5.29 ng/ul	35022	1,4-Dichlorobenzene-d4	8.048

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	56
2			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	56
3			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
4			3-Hexanol, 2,4-dimethyl-	130	C8H18O	013432-25-2	25
5			3-Pentanol, 2,4-dimethyl-	116	C7H16O	000600-36-2	17



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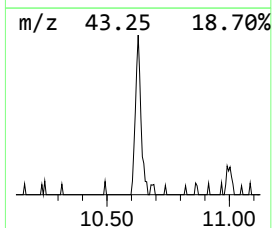
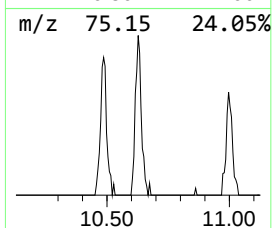
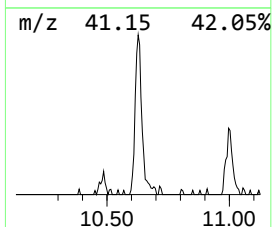
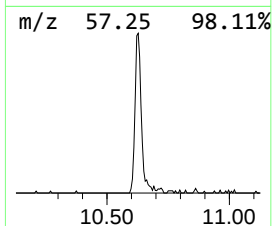
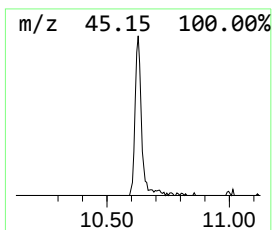
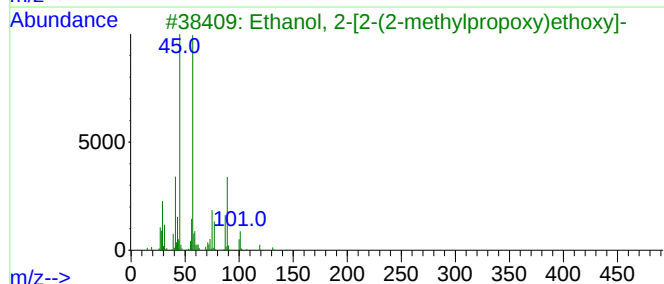
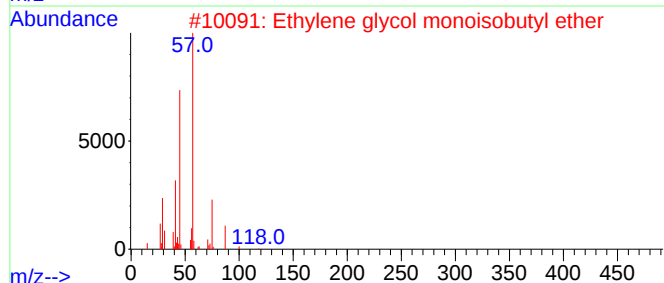
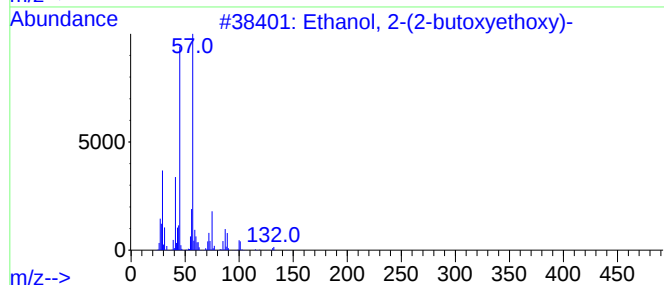
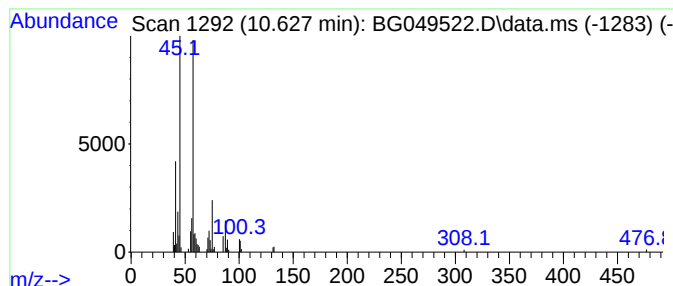
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG072421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.627	12.73 ng/ul	123996	Naphthalene-d8	10.868

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	86
2			Ethylene glycol monoisobutyl ether	118	C6H14O2	004439-24-1	72
3			Ethanol, 2-[2-(2-methylpropoxy)ethoxy]-	162	C8H18O3	018912-80-6	72
4			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	72
5			(2-Ethoxyethoxy)acetic acid, TBD...	262	C12H26O4Si	1000366-92-5	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072621\
Data File : BG049522.D
Acq On : 28 Jul 2021 00:51
Operator : CG/JU
Sample : M3091-01
Misc :
ALS Vial : 52 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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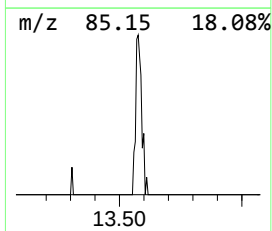
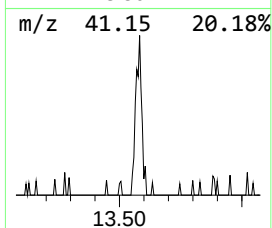
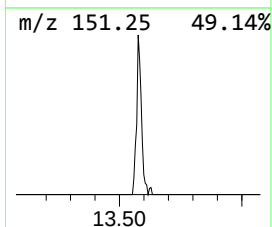
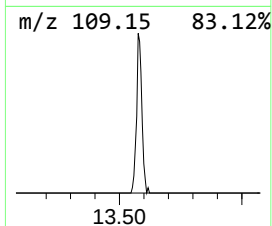
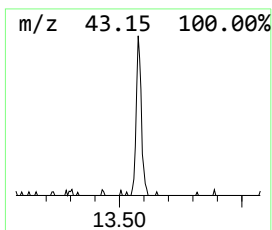
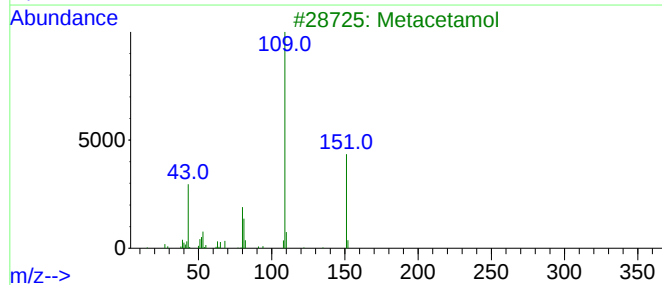
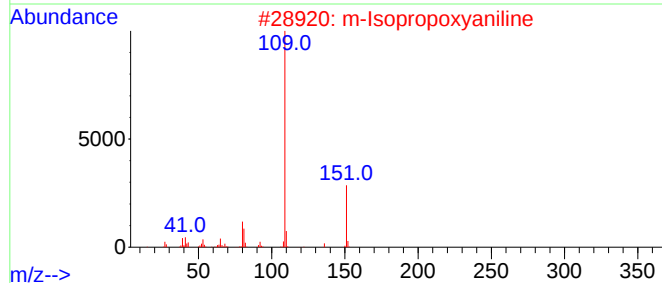
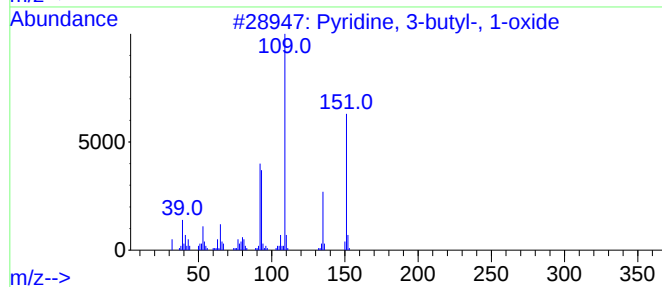
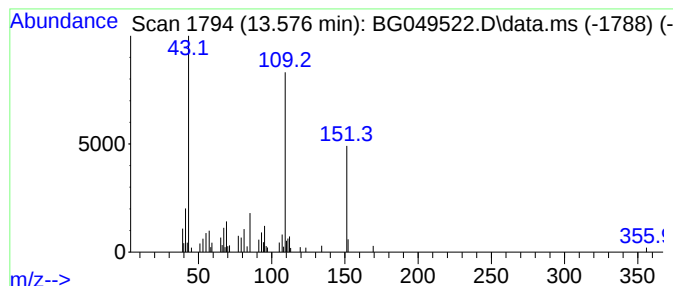
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Pyridine, 3-butyl-, 1-oxide Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.576	4.19 ng/ul	63511	Acenaphthene-d10	14.698

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pyridine, 3-butyl-, 1-oxide	151	C9H13NO	031396-33-5	64
2			m-Isopropoxyaniline	151	C9H13NO	041406-00-2	64
3			Metacetamol	151	C8H9NO2	000621-42-1	59
4			Metacetamol	151	C8H9NO2	000621-42-1	59
5			Metacetamol	151	C8H9NO2	000621-42-1	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072621\
Data File : BG049522.D
Acq On : 28 Jul 2021 00:51
Operator : CG/JU
Sample : M3091-01
Misc :
ALS Vial : 52 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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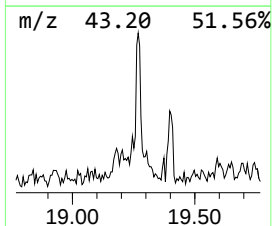
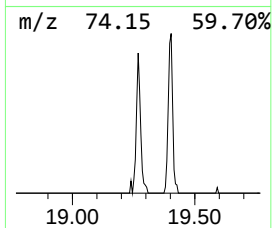
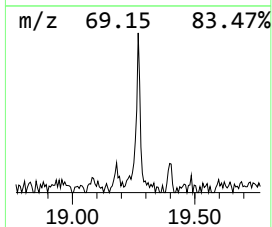
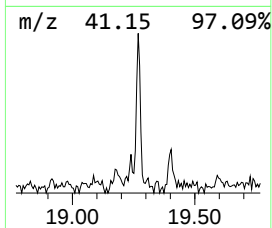
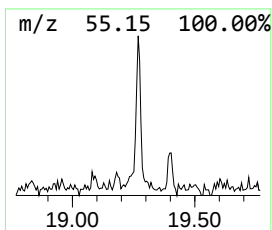
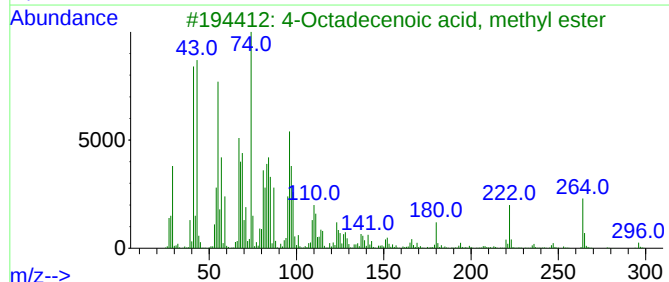
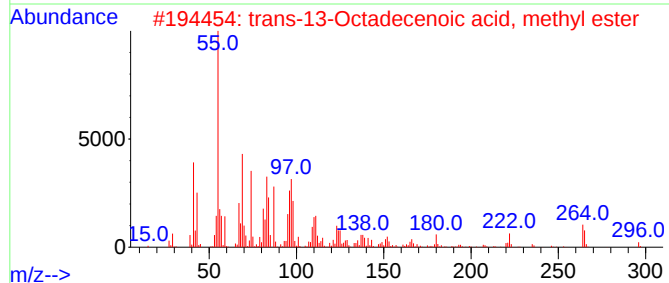
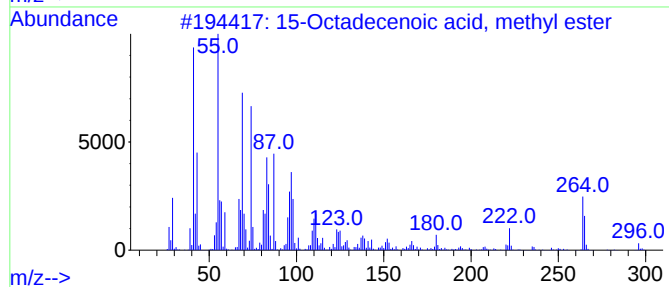
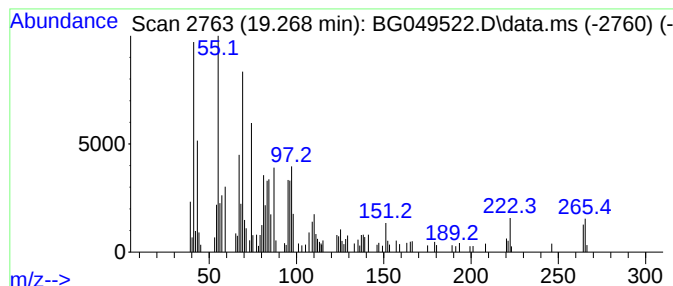
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 15-Octadecenoic acid, methyl... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.268	5.64 ng/ul	104893	Phenanthrene-d10	17.447

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			15-Octadecenoic acid, methyl ester	296	C19H36O2	004764-72-1	52
2			trans-13-Octadecenoic acid, methyl ester	296	C19H36O2	1000333-61-3	50
3			4-Octadecenoic acid, methyl ester	296	C19H36O2	056555-10-3	43
4			Trichloroacetic acid, undec-10-e...	314	C13H21Cl3O2	1000280-51-3	43
5			2,6,10-Dodecatrienal, 3,7,11-tri...	220	C15H24O	004380-32-9	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072621\
Data File : BG049522.D
Acq On : 28 Jul 2021 00:51
Operator : CG/JU
Sample : M3091-01
Misc :
ALS Vial : 52 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0036

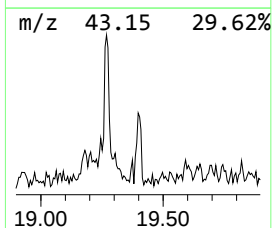
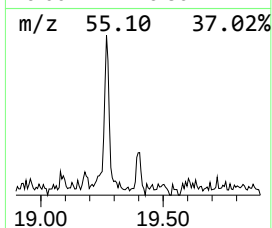
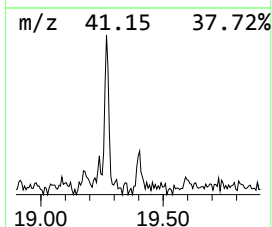
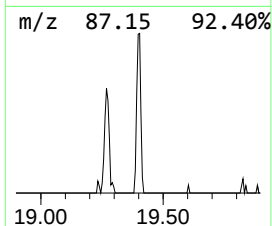
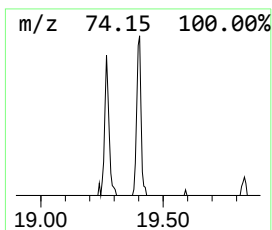
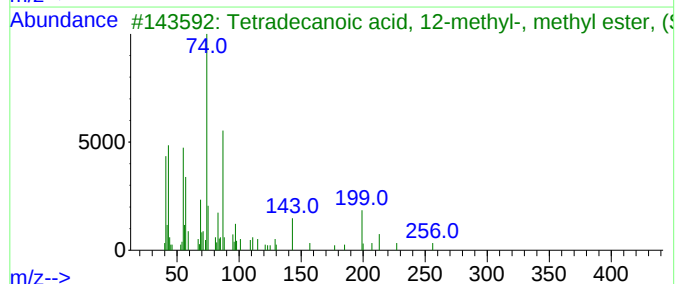
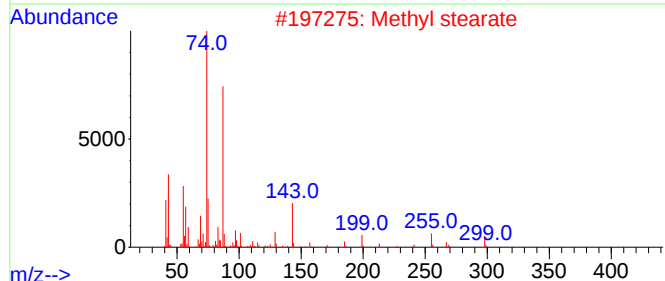
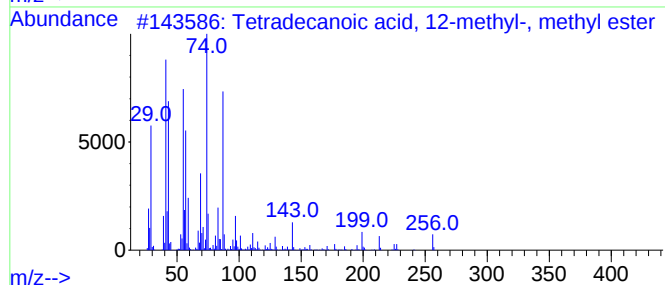
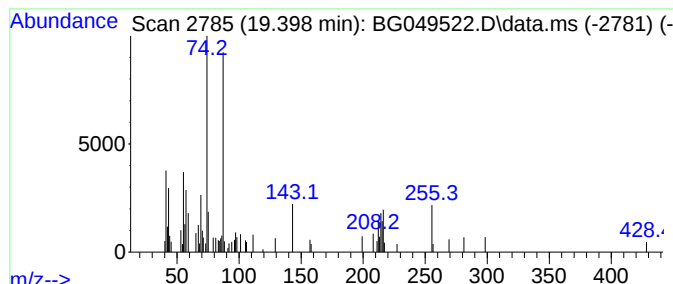
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Tetradecanoic acid, 12-meth... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.398	2.12 ng/ul	39389	Phenanthrene-d10	17.447

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetradecanoic acid, 12-methyl-, ...	256	C16H32O2	005129-66-8	70
2			Methyl stearate	298	C19H38O2	000112-61-8	68
3			Tetradecanoic acid, 12-methyl-, ...	256	C16H32O2	062691-05-8	64
4			Pentadecanoic acid, methyl ester	256	C16H32O2	007132-64-1	64
5			Methyl 9-methyltetradecanoate	256	C16H32O2	213617-69-7	62



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072621\
Data File : BG049522.D
Acq On : 28 Jul 2021 00:51
Operator : CG/JU
Sample : M3091-01
Misc :
ALS Vial : 52 Sample Multiplier: 1

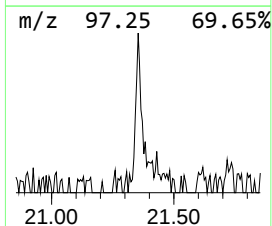
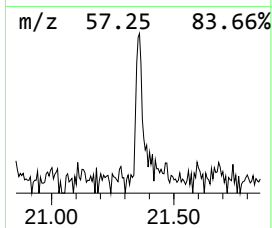
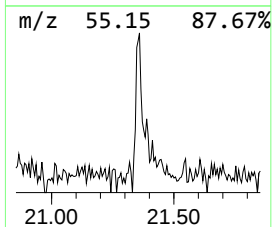
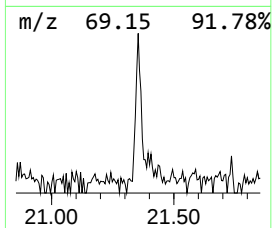
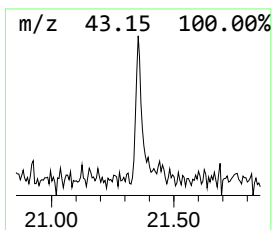
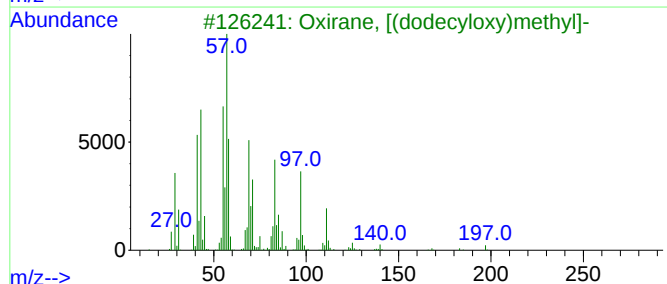
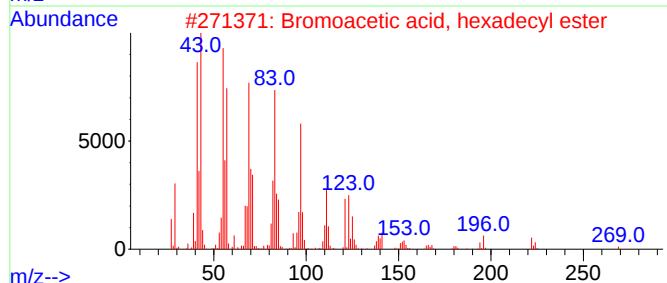
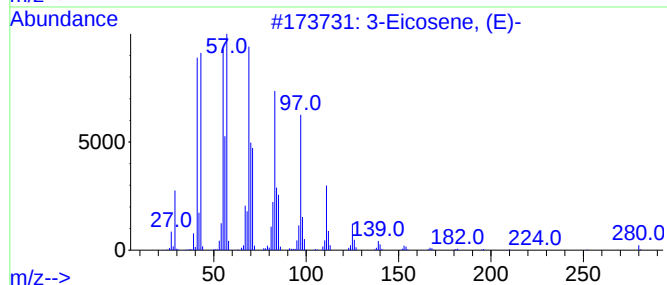
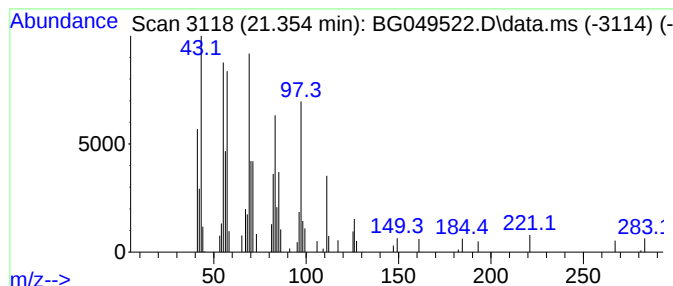
Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG072421.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD		R.T.	
21.354	2.76 ng/ul	52074	Chrysene-d12		21.730	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	3-Eicosene, (E)-		280	C20H40	074685-33-9	91
2	Bromoacetic acid, hexadecyl ester		362	C18H35BrO2	005454-48-8	91
3	Oxirane, [(dodecyloxy)methyl]-		242	C15H30O2	002461-18-9	80
4	Trichloroacetic acid, pentadecyl...		372	C17H31Cl3O2	074339-53-0	68
5	Cyclotetradecane		196	C14H28	000295-17-0	68



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072621\
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Acq On : 28 Jul 2021 00:51
Operator : CG/JU
Sample : M3091-01
Misc :
ALS Vial : 52 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG072421.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
(DEL) Alkane: C...	3.078	11.4	ng/ul	75427	1	8.048	132441	20.0
Butane, 2-metho...	3.161	5.3	ng/ul	35022	1	8.048	132441	20.0
Ethanol, 2-(2-b...	10.627	12.7	ng/ul	123996	2	10.868	194762	20.0
Pyridine, 3-but...	13.576	4.2	ng/ul	63511	3	14.698	303346	20.0
15-Octadecenoic...	19.268	5.6	ng/ul	104893	4	17.447	371943	20.0
Tetradecanoic a...	19.398	2.1	ng/ul	39389	4	17.447	371943	20.0
(DEL) Alkane: S...	21.354	2.8	ng/ul	52074	5	21.730	377255	20.0