

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
 Data File : BG049651.D
 Acq On : 5 Aug 2021 15:55
 Operator : CG/JU
 Sample : M3162-06
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BG06

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-BG080421.M
 Title : SVOA CALIBRATION

Signal : TIC: BG049651.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.077	3	6	14	rVB2	44574	79779	9.23%	0.863%
2	3.153	15	19	30	rVB	41186	76116	8.80%	0.824%
3	3.229	30	32	34	rBV3	2246	2037	0.24%	0.022%
4	3.329	46	49	50	rBV3	1787	2036	0.24%	0.022%
5	3.429	64	66	72	rVB5	6838	9204	1.06%	0.100%
6	3.476	72	74	78	rBV3	1870	1894	0.22%	0.020%
7	3.846	134	137	149	rBV	58547	120151	13.89%	1.300%
8	4.152	186	189	192	rBV2	1607	2073	0.24%	0.022%
9	4.181	192	194	200	rVB5	2532	4385	0.51%	0.047%
10	4.410	229	233	235	rBV3	2511	2532	0.29%	0.027%
11	5.127	353	355	358	rVB4	1820	1906	0.22%	0.021%
12	5.879	479	483	486	rBV3	1199	1856	0.21%	0.020%
13	6.942	659	664	665	rBV2	1429	1735	0.20%	0.019%
14	7.195	701	707	718	rBV	120926	224855	26.00%	2.433%
15	7.359	728	735	744	rVB	75854	129110	14.93%	1.397%
16	7.571	764	771	779	rBV	110083	198982	23.01%	2.153%
17	8.040	845	851	857	rBV2	89348	158181	18.29%	1.711%
18	8.745	964	971	979	rBV2	148880	266095	30.77%	2.879%
19	9.209	1043	1050	1060	rVB	120398	217209	25.12%	2.350%
20	9.374	1076	1078	1080	rBV2	2215	1806	0.21%	0.020%
21	9.621	1116	1120	1122	rBV3	1866	2945	0.34%	0.032%
22	9.938	1166	1174	1185	rVB2	107519	203488	23.53%	2.202%
23	10.020	1185	1188	1191	rBV	1170	1691	0.20%	0.018%
24	10.478	1258	1266	1278	rVB2	150935	278838	32.24%	3.017%
25	10.619	1284	1290	1304	rBV3	43394	84728	9.80%	0.917%
26	10.860	1325	1331	1339	rVB	121550	221875	25.66%	2.401%
27	10.995	1347	1354	1363	rBV2	146868	266846	30.86%	2.887%
28	11.900	1504	1508	1516	rBV2	11364	23280	2.69%	0.252%
29	12.446	1599	1601	1605	rVB2	2183	1920	0.22%	0.021%
30	13.504	1777	1781	1786	rBV3	2663	3275	0.38%	0.035%
31	13.574	1786	1793	1799	rVB2	11460	17413	2.01%	0.188%
32	13.674	1804	1810	1813	rVB2	1163	1862	0.22%	0.020%
33	14.091	1874	1881	1888	rBV	288381	426951	49.37%	4.619%
34	14.385	1924	1931	1938	rBV	311504	465478	53.83%	5.036%
35	14.690	1976	1983	1991	rVB	200041	309358	35.77%	3.347%
36	14.884	2009	2016	2028	rBV	109080	199446	23.06%	2.158%

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Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

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Start Thrs: 0.2

Max Peaks: 100

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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-BG080421.M
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37	15.683	2146	2152	2162	rVB	376119	575760	66.58%	6.229%
38	15.801	2166	2172	2179	rBV	135915	200016	23.13%	2.164%
39	15.924	2187	2193	2201	rVB4	3772	7337	0.85%	0.079%
40	16.130	2225	2228	2232	rBV	1704	2859	0.33%	0.031%
41	17.439	2444	2451	2458	rBV	242175	374195	43.27%	4.049%
42	17.539	2462	2468	2476	rVB	389251	610069	70.55%	6.601%
43	17.774	2505	2508	2512	rBV	1718	2388	0.28%	0.026%
44	18.097	2560	2563	2568	rBV3	12420	13683	1.58%	0.148%
45	18.256	2588	2590	2593	rBV	1651	1842	0.21%	0.020%
46	18.321	2595	2601	2606	rBV3	8643	13709	1.59%	0.148%
47	18.755	2671	2675	2685	rBV	36359	59111	6.84%	0.640%
48	19.114	2733	2736	2737	rVB	2138	1896	0.22%	0.021%
49	19.143	2737	2741	2744	rBV2	1825	3132	0.36%	0.034%
50	19.231	2754	2756	2757	rBV	2519	1744	0.20%	0.019%
51	19.266	2758	2762	2766	rVV2	34247	40285	4.66%	0.436%
52	19.331	2769	2773	2775	rBV2	1831	2924	0.34%	0.032%
53	19.402	2780	2785	2791	rVB2	11962	16026	1.85%	0.173%
54	19.466	2793	2796	2799	rBV2	4534	6507	0.75%	0.070%
55	19.543	2806	2809	2811	rBV2	1503	1884	0.22%	0.020%
56	19.830	2852	2858	2867	rBV	466791	691208	79.93%	7.478%
57	20.124	2906	2908	2910	rBV3	2147	1738	0.20%	0.019%
58	20.159	2912	2914	2916	rBV3	1779	2093	0.24%	0.023%
59	20.330	2941	2943	2944	rBV	2127	1805	0.21%	0.020%
60	20.347	2944	2946	2954	rVB4	4802	8068	0.93%	0.087%
61	20.447	2961	2963	2966	rVB2	2468	2455	0.28%	0.027%
62	20.524	2975	2976	2979	rVB2	3239	2379	0.28%	0.026%
63	20.582	2984	2986	2991	rVV4	2662	3325	0.38%	0.036%
64	20.653	2995	2998	3001	rBV4	1974	2652	0.31%	0.029%
65	20.676	3001	3002	3006	rVB3	2993	2562	0.30%	0.028%
66	20.729	3008	3011	3012	rBV2	4069	4821	0.56%	0.052%
67	20.800	3018	3023	3028	rBV	782383	864780	100.00%	9.356%
68	20.841	3028	3030	3031	rVV2	4294	2731	0.32%	0.030%
69	20.853	3031	3032	3036	rVB4	2666	3169	0.37%	0.034%
70	21.352	3113	3117	3125	rBV4	22188	39654	4.59%	0.429%
71	21.487	3138	3140	3142	rBV3	2639	2696	0.31%	0.029%
72	21.599	3155	3159	3164	rBV4	6185	10708	1.24%	0.116%
73	21.640	3164	3166	3168	rBV3	2254	2629	0.30%	0.028%
74	21.722	3173	3180	3187	rVV	235095	391019	45.22%	4.231%
75	22.157	3250	3254	3258	rVB5	3498	4955	0.57%	0.054%
76	22.380	3290	3292	3295	rBV4	2385	1742	0.20%	0.019%

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 ClientSampleId :
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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

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77	22.515	3310	3315	3316	rBV5	6984	11425	1.32%	0.124%
78	22.527	3316	3317	3324	rVB6	11345	11564	1.34%	0.125%
79	23.061	3406	3408	3410	rBV3	2588	2751	0.32%	0.030%
80	23.132	3418	3420	3422	rBV3	3070	2377	0.27%	0.026%
81	23.901	3546	3551	3555	rVB7	10159	18796	2.17%	0.203%
82	24.042	3571	3575	3582	rVB5	13836	29972	3.47%	0.324%
83	24.242	3607	3609	3611	rBV3	2518	2015	0.23%	0.022%
84	24.771	3689	3699	3709	rBV2	239882	656670	75.93%	7.105%
85	25.000	3727	3738	3748	rVB2	142211	422123	48.81%	4.567%
86	25.940	3896	3898	3899	rVB2	3248	1957	0.23%	0.021%
87	25.957	3899	3901	3905	rVB4	3625	3269	0.38%	0.035%
88	25.993	3905	3907	3910	rBV4	2739	3585	0.41%	0.039%
89	26.169	3930	3937	3948	rVB7	18459	54905	6.35%	0.594%
90	26.374	3970	3972	3973	rVB	3559	2162	0.25%	0.023%
91	26.674	4021	4023	4024	rBV2	3098	1896	0.22%	0.021%
92	27.890	4228	4230	4231	rBV	2426	1744	0.20%	0.019%
93	28.918	4403	4405	4407	rVB2	2721	2105	0.24%	0.023%
94	29.171	4447	4448	4450	rBV2	2676	2443	0.28%	0.026%
95	30.075	4600	4602	4604	rVB3	3374	2430	0.28%	0.026%
96	30.234	4627	4629	4630	rBV	3233	2505	0.29%	0.027%
97	30.404	4655	4658	4661	rBV4	4005	5041	0.58%	0.055%
98	31.197	4792	4793	4794	rBV	3934	2233	0.26%	0.024%
99	31.491	4841	4843	4845	rBV3	2849	1928	0.22%	0.021%
100	31.661	4871	4872	4874	rBV3	3436	2900	0.34%	0.031%

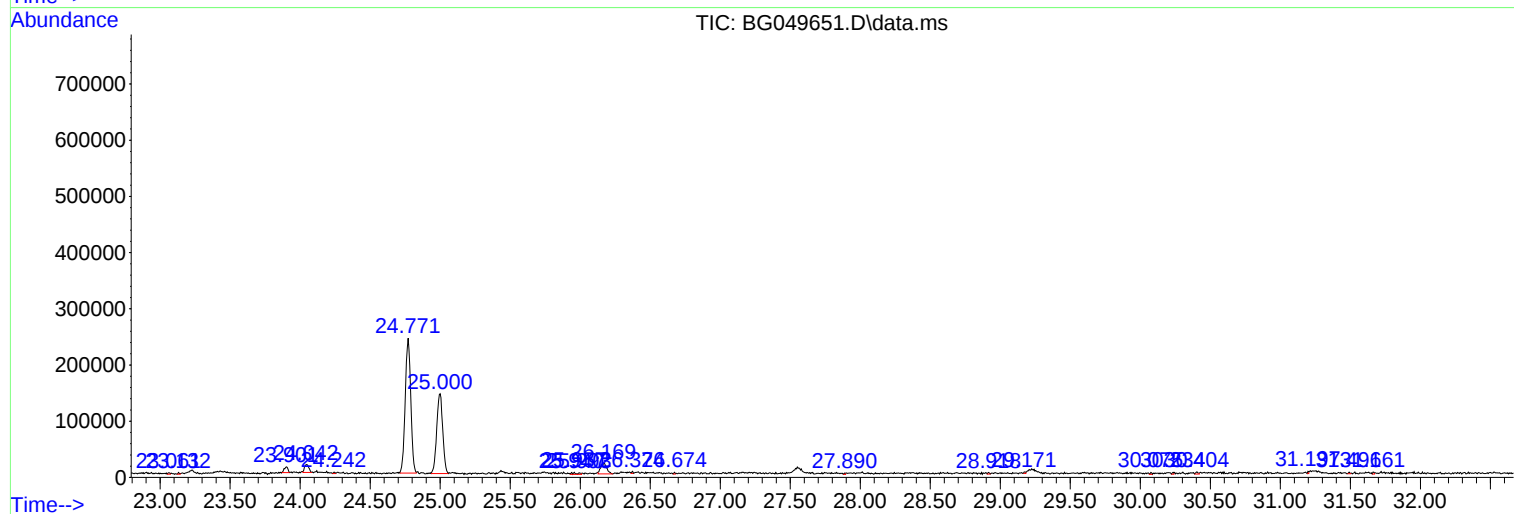
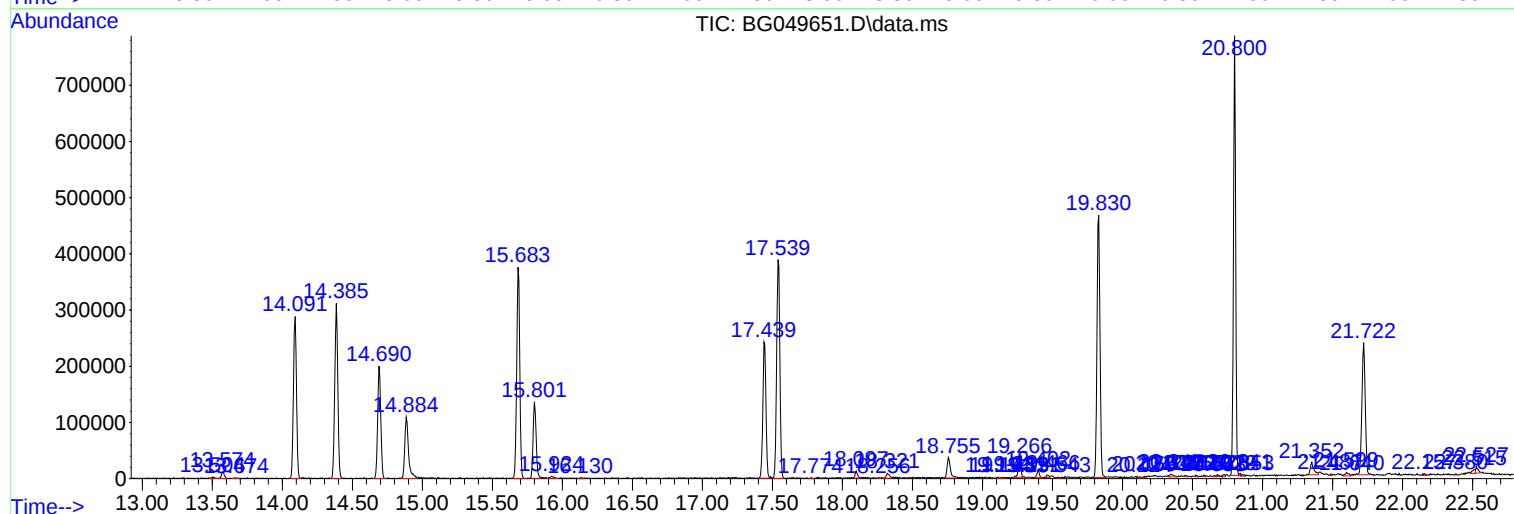
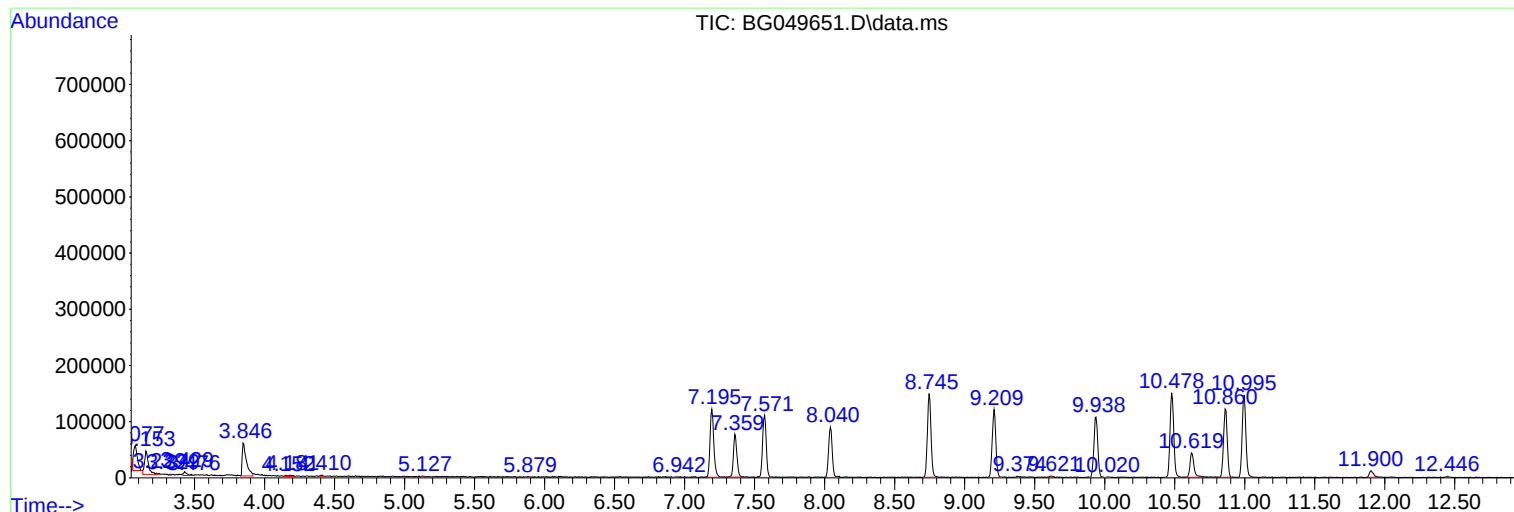
Sum of corrected areas: 9242688

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Instrument :
BNA_G
ClientSampleId :
BGET6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
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Acq On : 5 Aug 2021 15:55
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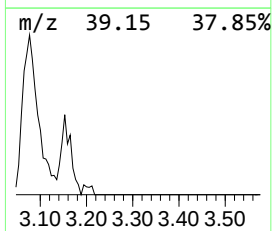
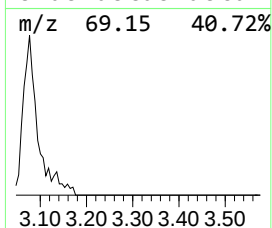
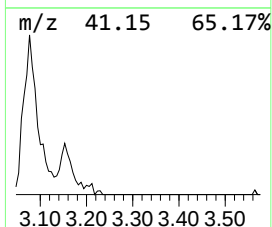
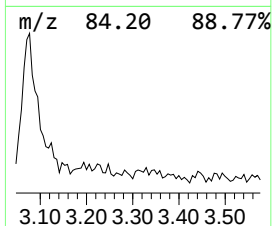
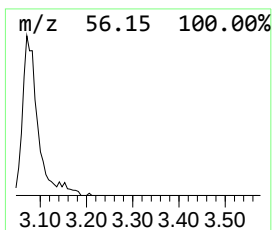
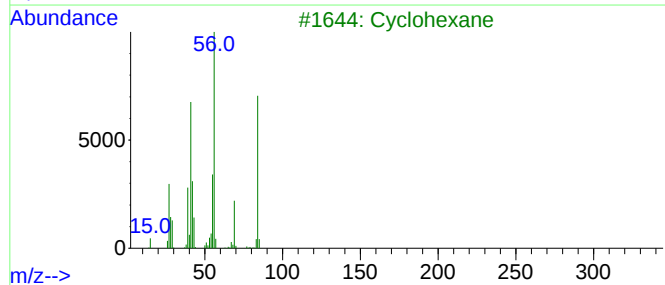
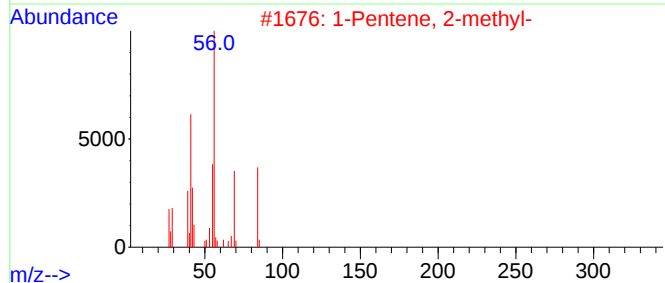
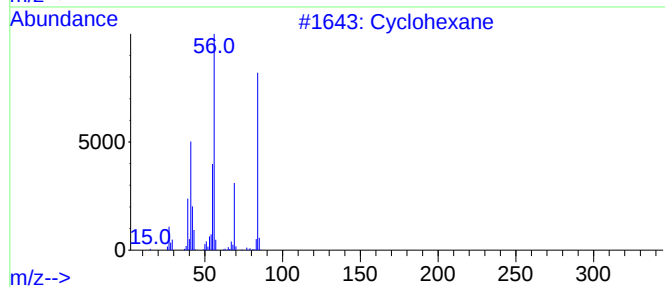
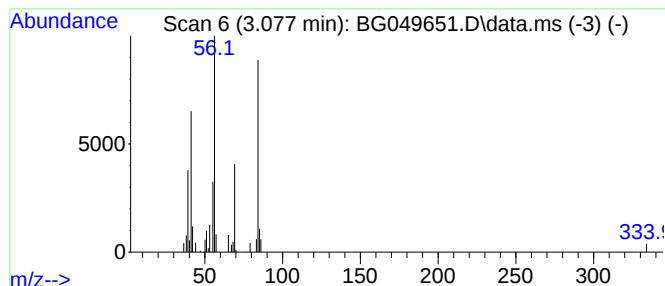
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 (DEL) Alkane: Cyclic3.077 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.077	10.09 ng/ul	79779	1,4-Dichlorobenzene-d4	8.040

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclohexane	84	C6H12	000110-82-7	72
2		1-Pentene, 2-methyl-	84	C6H12	000763-29-1	64
3		Cyclohexane	84	C6H12	000110-82-7	64
4		Cyclohexane	84	C6H12	000110-82-7	64
5		Pyridine-D5-	84	C5D5N	007291-22-7	59



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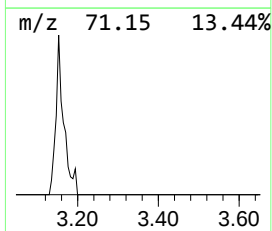
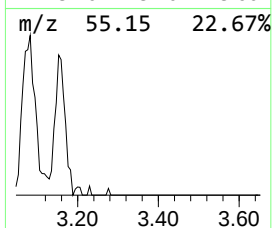
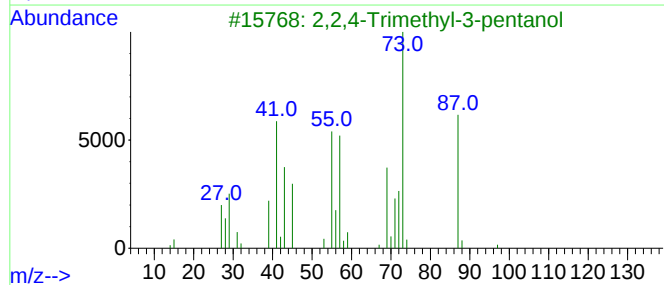
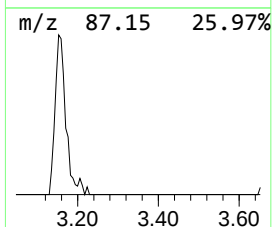
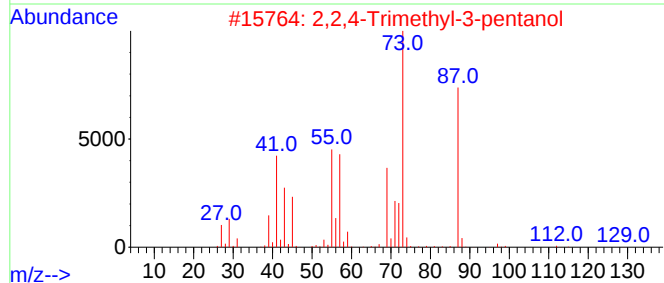
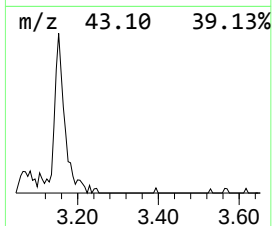
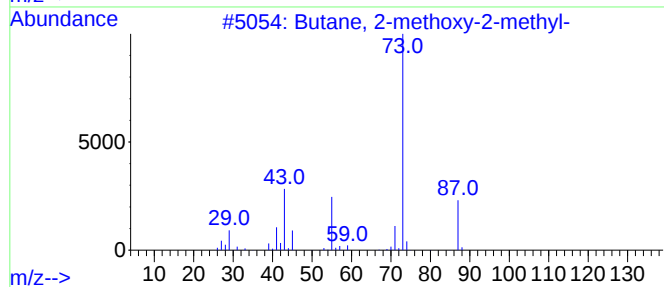
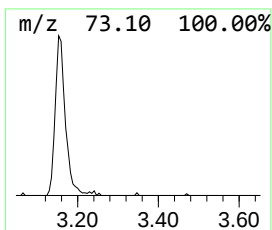
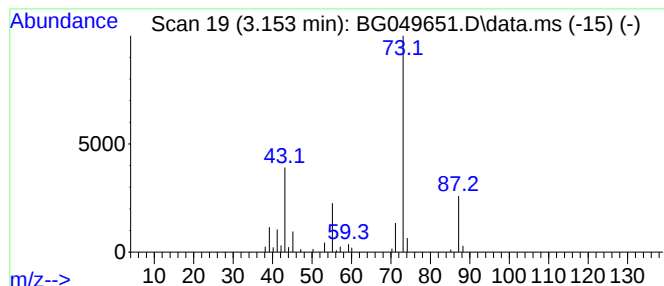
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Butane, 2-methoxy-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.153	9.62 ng/ul	76116	1,4-Dichlorobenzene-d4	8.040

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	56
3			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
4			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	40
5			N-Ethylformamide	73	C3H7NO	000627-45-2	9



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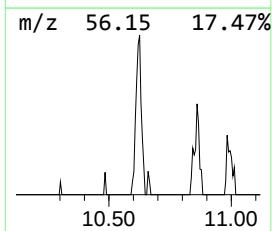
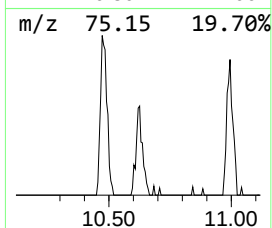
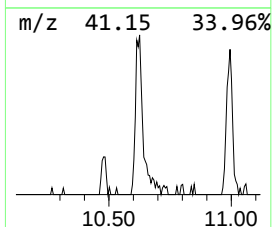
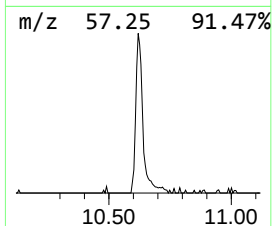
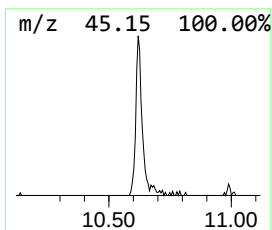
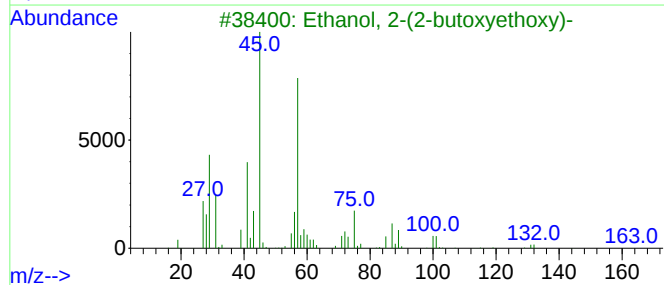
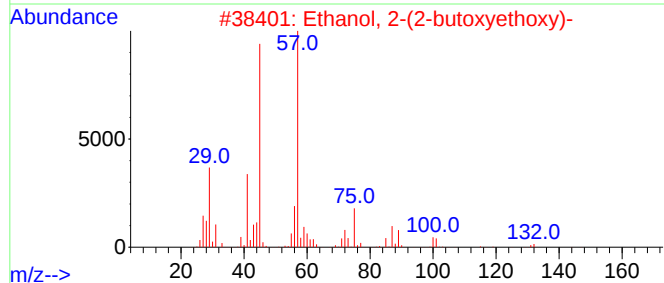
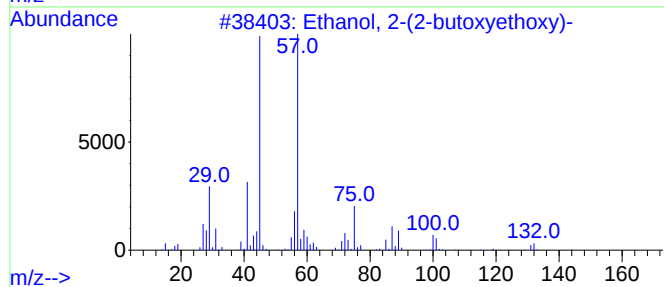
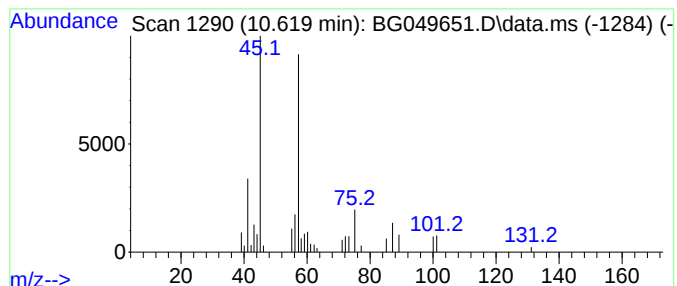
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.619	7.64 ng/ul	84728	Naphthalene-d8	10.860

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	86
3			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	78
4			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	72
5			Propanoic acid, 2-hydroxy-, 2-me...	146	C7H14O3	000585-24-0	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049651.D
Acq On : 5 Aug 2021 15:55
Operator : CG/JU
Sample : M3162-06
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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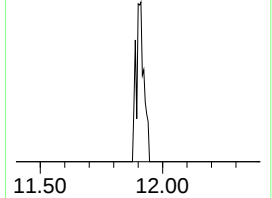
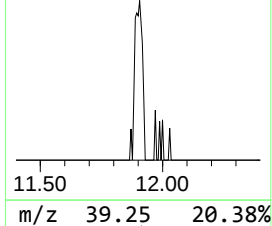
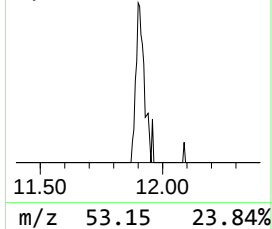
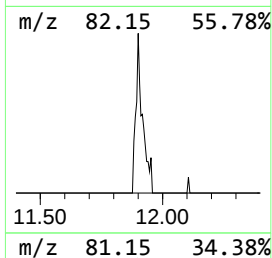
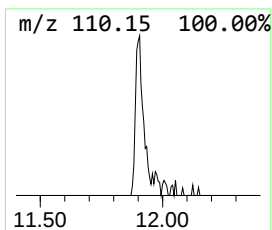
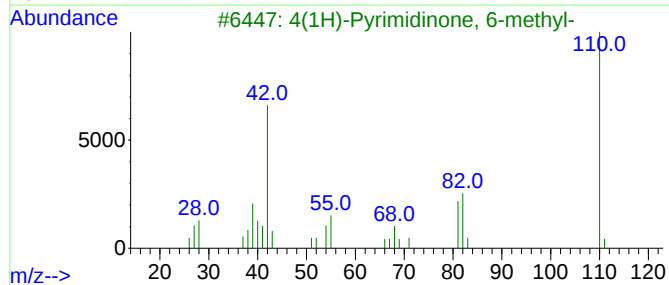
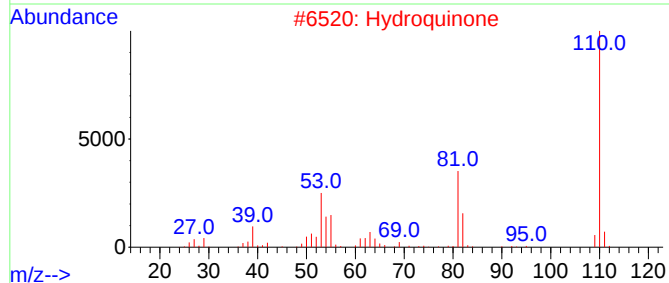
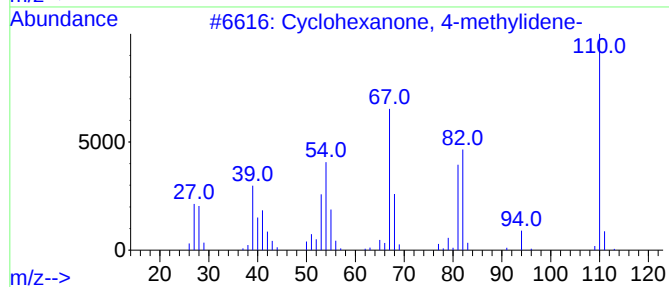
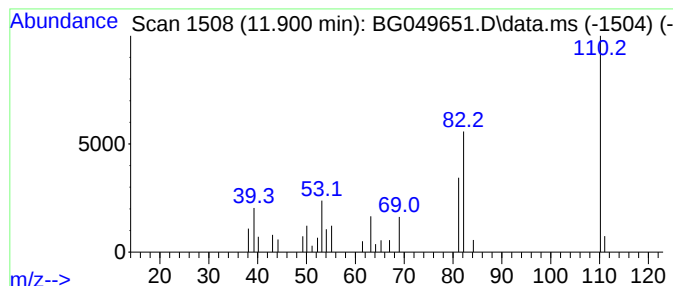
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Cyclohexanone, 4-methylidene- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.900	2.10 ng/ul	23280	Naphthalene-d8	10.860

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexanone, 4-methylidene-	110	C7H10O	029648-66-6	59
2			Hydroquinone	110	C6H6O2	000123-31-9	58
3			4(1H)-Pyrimidinone, 6-methyl-	110	C5H6N2O	003524-87-6	58
4			1-Methylimidazole-4-carboxaldehyde	110	C5H6N2O	017289-26-8	53
5			4(1H)-Pyrimidinone, 1-methyl-	110	C5H6N2O	002228-30-0	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049651.D
Acq On : 5 Aug 2021 15:55
Operator : CG/JU
Sample : M3162-06
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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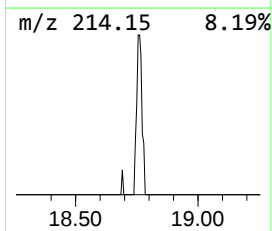
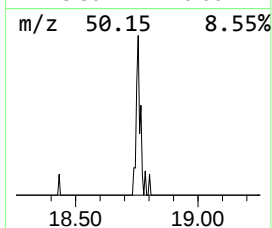
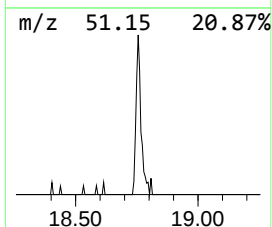
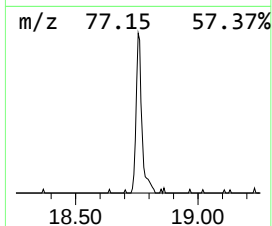
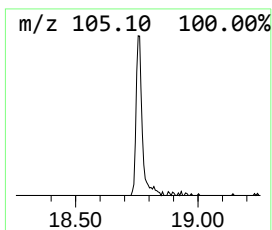
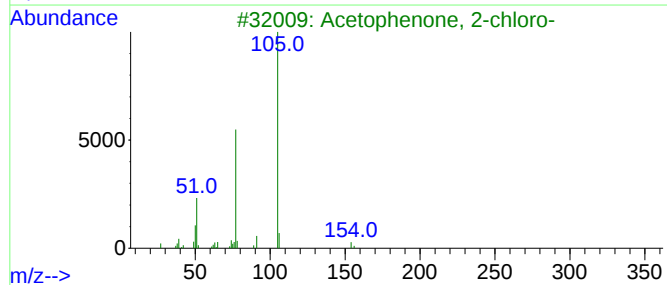
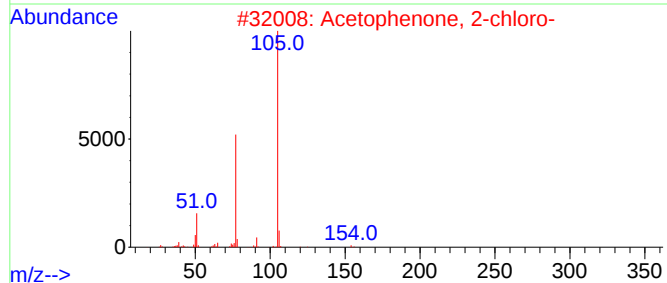
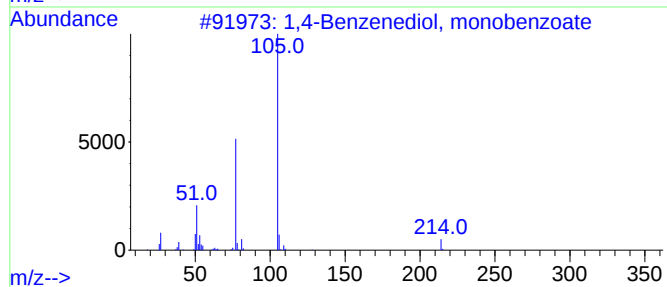
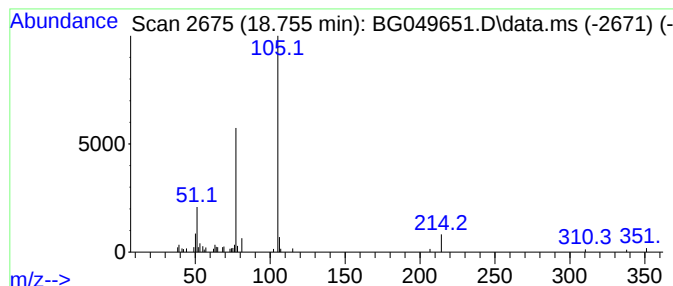
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 1,4-Benzenediol, monobenzoate Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.755	3.16 ng/ul	59111	Phenanthrene-d10	17.445

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,4-Benzenediol, monobenzoate	214	C13H10O3	002444-19-1	80
2			Acetophenone, 2-chloro-	154	C8H7ClO	000532-27-4	72
3			Acetophenone, 2-chloro-	154	C8H7ClO	000532-27-4	72
4			.beta.-Benzilmonoxime	225	C14H11NO2	014090-77-8	72
5			Glyoxylic acid, phenyl-, 2'-cyan...	251	C15H9NO3	166538-14-3	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049651.D
Acq On : 5 Aug 2021 15:55
Operator : CG/JU
Sample : M3162-06
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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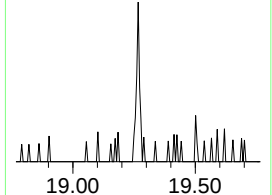
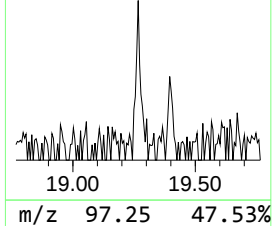
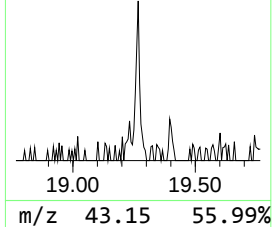
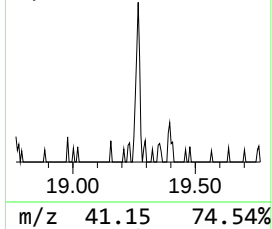
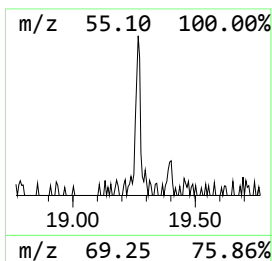
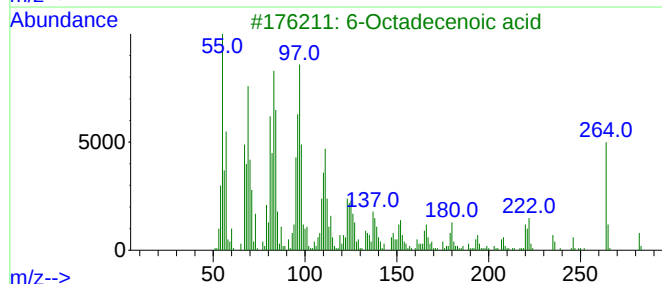
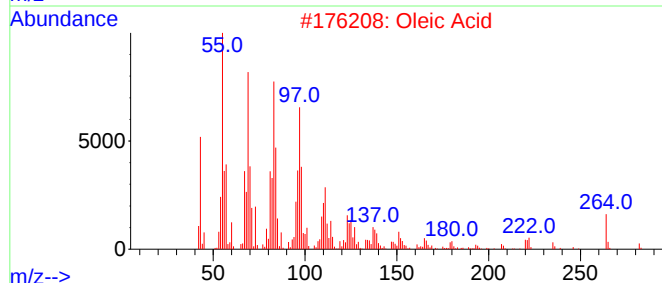
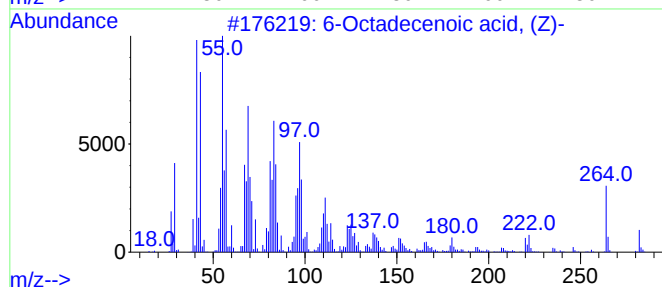
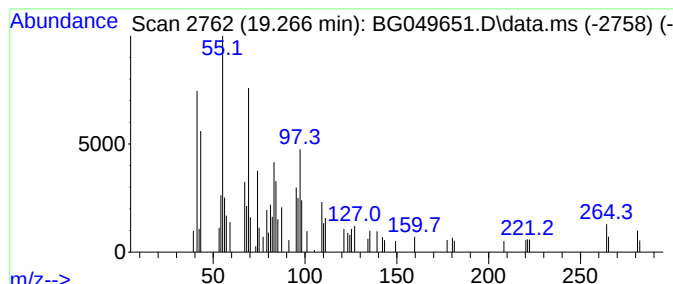
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 6-Octadecenoic acid, (Z)- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.266	2.15 ng/ul	40285	Phenanthrene-d10	17.445

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			6-Octadecenoic acid, (Z)-	282	C18H34O2	000593-39-5	95
2			Oleic Acid	282	C18H34O2	000112-80-1	89
3			6-Octadecenoic acid	282	C18H34O2	1000336-66-8	87
4			9-Octadecenoic acid, (E)-	282	C18H34O2	000112-79-8	83
5			Cyclopentanone, 3-[3,5-decadieny...	220	C15H24O	1000131-99-7	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049651.D
Acq On : 5 Aug 2021 15:55
Operator : CG/JU
Sample : M3162-06
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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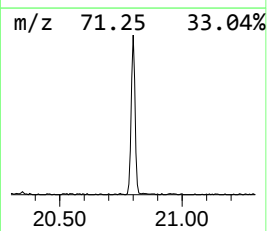
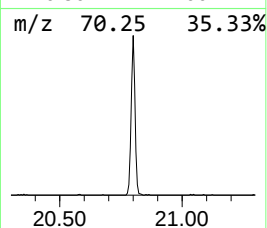
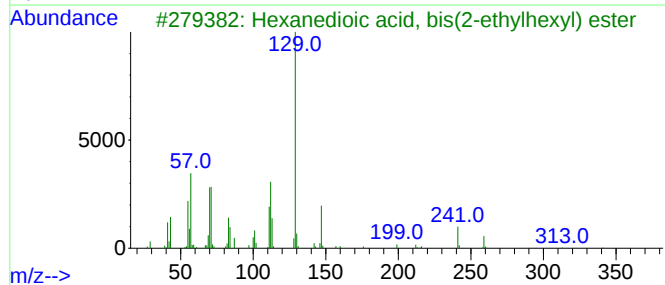
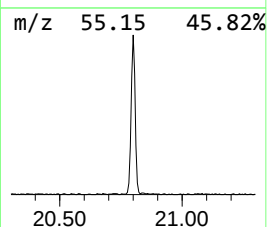
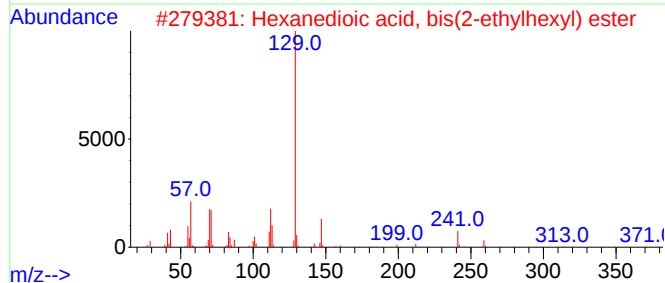
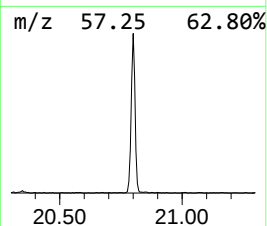
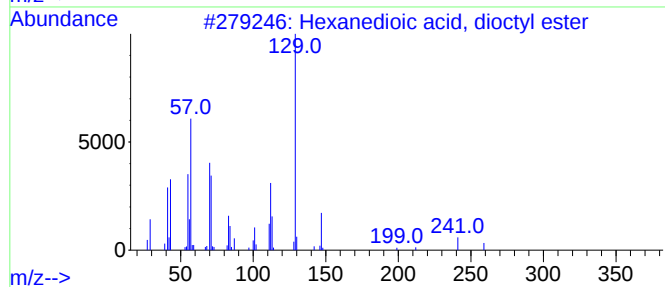
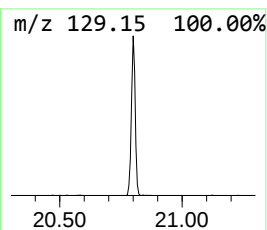
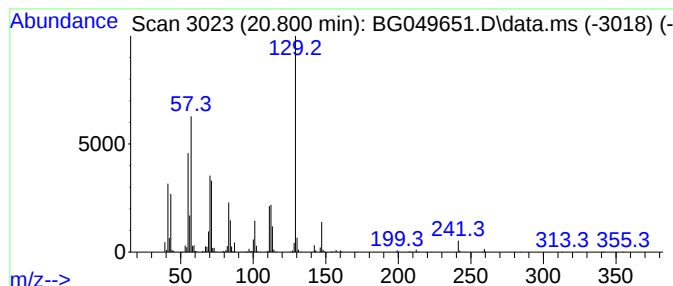
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Hexanedioic acid, dioctyl e... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.800	44.23 ng/ul	864780	Chrysene-d12	21.722

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexanedioic acid, dioctyl ester	370	C22H42O4	000123-79-5	87
2			Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	80
3			Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	80
4			Diisooctyl adipate	370	C22H42O4	001330-86-5	78
5			Adipic acid, octyl 2-octyl ester	370	C22H42O4	1000324-45-0	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049651.D
Acq On : 5 Aug 2021 15:55
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Sample : M3162-06
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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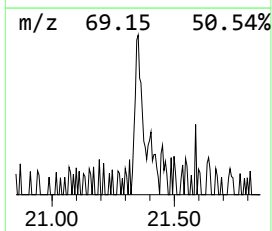
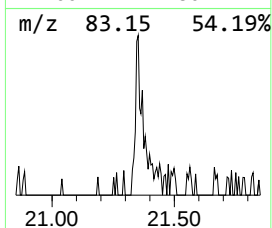
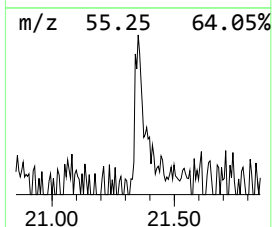
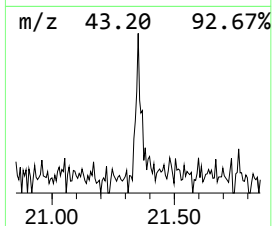
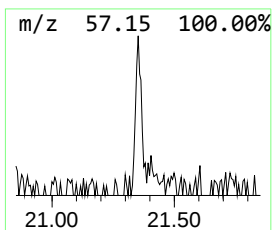
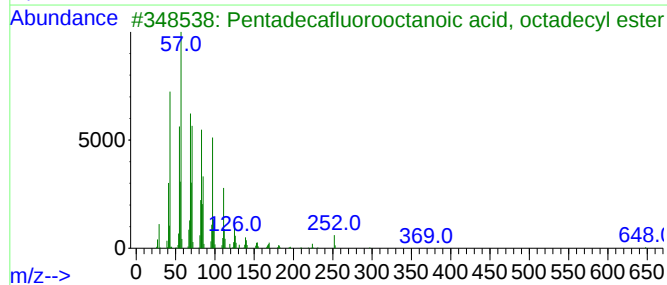
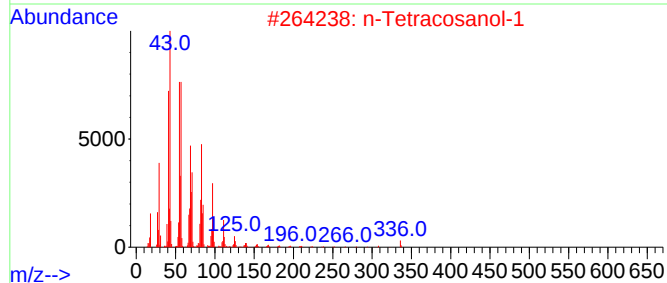
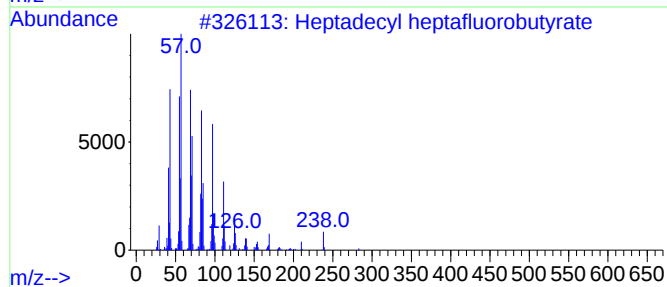
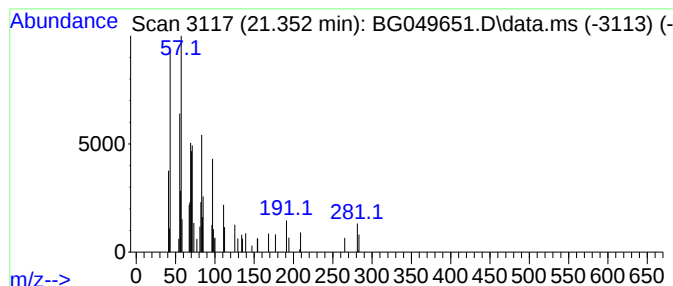
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Heptadecyl heptafluorobutyrate Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.352	2.03 ng/ul	39654	Chrysene-d12	21.722

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	76
2			n-Tetracosanol-1	354	C24H50O	000506-51-4	76
3			Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	76
4			17-Pentatriacontene	491	C35H70	006971-40-0	70
5			Octadecyl trifluoroacetate	366	C20H37F3O2	079392-43-1	68



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049651.D
Acq On : 5 Aug 2021 15:55
Operator : CG/JU
Sample : M3162-06
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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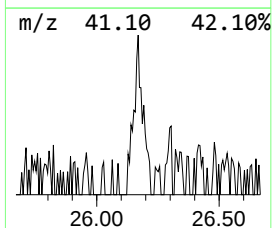
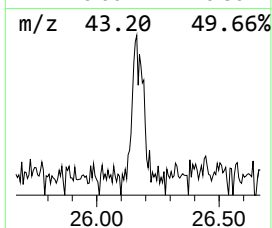
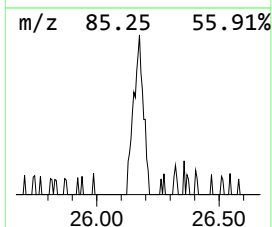
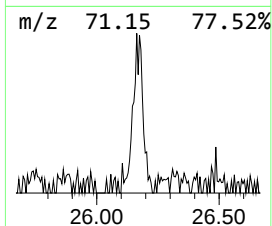
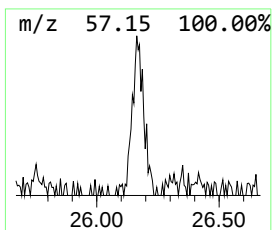
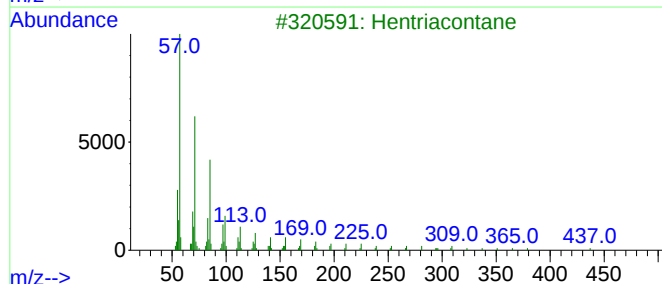
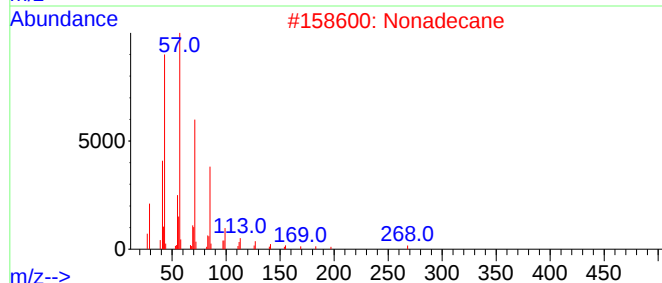
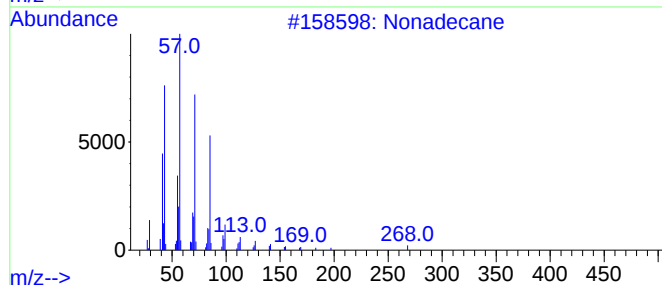
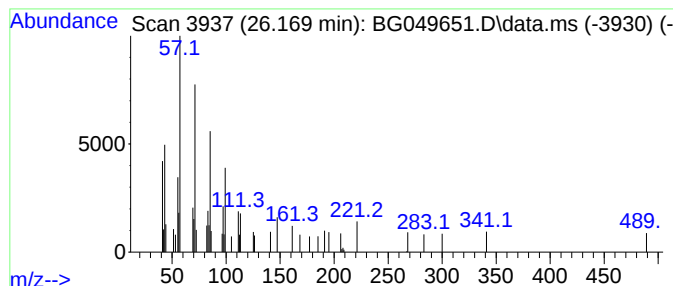
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.169	2.60 ng/ul	54905	Perylene-d12	24.994

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonadecane	268	C19H40	000629-92-5	83
2			Nonadecane	268	C19H40	000629-92-5	58
3			Hentriacontane	437	C31H64	000630-04-6	58
4			Dotriacontane, 1-iodo-	576	C32H65I	1000406-32-4	53
5			Tetrapentacontane	759	C54H110	005856-66-6	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049651.D
Acq On : 5 Aug 2021 15:55
Operator : CG/JU
Sample : M3162-06
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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Butane, 2-metho...	3.153	9.6	ng/ul	76116	1	8.040	158181	20.0
Ethanol, 2-(2-b...	10.619	7.6	ng/ul	84728	2	10.860	221875	20.0
Cyclohexanone, ...	11.900	2.1	ng/ul	23280	2	10.860	221875	20.0
1,4-Benzenediol...	18.755	3.2	ng/ul	59111	4	17.445	374195	20.0
6-Octadecenoic ...	19.266	2.1	ng/ul	40285	4	17.445	374195	20.0
Hexanedioic aci...	20.800	44.2	ng/ul	864780	5	21.722	391019	20.0
Heptadecyl hept...	21.352	2.0	ng/ul	39654	5	21.722	391019	20.0
(DEL) Alkane: S...	26.169	2.6	ng/ul	54905	6	24.994	422123	20.0