

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021921\
 Data File : BG048225.D
 Acq On : 20 Feb 2021 6:13
 Operator : CG/JU
 Sample : M1412-02
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 DBG5

Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : OFF
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 0.1 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG021321MA.M
 Title : SVOA CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.509	25	29	37	rVB5	6748	11521	1.15%	0.091%
2	4.043	116	120	123	rVB4	2520	4265	0.43%	0.034%
3	4.161	135	140	145	rBV3	4448	7600	0.76%	0.060%
4	4.273	155	159	167	rVB5	3012	5930	0.59%	0.047%
5	5.195	311	316	324	rVB2	8445	15951	1.59%	0.126%
6	5.248	324	325	328	rVB2	1850	1712	0.17%	0.014%
7	6.629	558	560	562	rVB2	2435	2040	0.20%	0.016%
8	6.729	574	577	580	rVB2	1359	1749	0.17%	0.014%
9	6.781	580	586	592	rBV	92321	143308	14.32%	1.130%
10	7.093	634	639	644	rBV3	6271	10150	1.01%	0.080%
11	7.287	659	672	684	rBV	107781	208522	20.84%	1.645%
12	7.428	689	696	707	rBV	79220	141185	14.11%	1.113%
13	7.516	707	711	712	rBV	5024	5461	0.55%	0.043%
14	7.645	727	733	742	rBV	113483	198342	19.82%	1.564%
15	7.845	763	767	770	rBV	1239	1712	0.17%	0.014%
16	8.109	803	812	819	rBV	125411	215233	21.51%	1.697%
17	8.239	828	834	841	rVB3	7722	12762	1.28%	0.101%
18	8.391	857	860	861	rBV3	1938	2050	0.20%	0.016%
19	8.409	861	863	868	rVB4	4346	6056	0.61%	0.048%
20	8.826	928	934	946	rBV2	117606	223007	22.28%	1.759%
21	8.944	950	954	957	rVB4	2645	3964	0.40%	0.031%
22	9.278	1003	1011	1022	rBV	107586	199491	19.93%	1.573%
23	9.419	1032	1035	1037	rBV2	1559	2093	0.21%	0.017%
24	9.666	1071	1077	1081	rVB4	2081	2801	0.28%	0.022%
25	9.766	1092	1094	1096	rBV2	1844	1840	0.18%	0.015%
26	10.001	1126	1134	1141	rBV	102083	180707	18.06%	1.425%
27	10.230	1167	1173	1177	rBV5	9634	15828	1.58%	0.125%
28	10.283	1179	1182	1184	rVB2	1767	2040	0.20%	0.016%
29	10.330	1186	1190	1193	rVB2	1923	2442	0.24%	0.019%
30	10.553	1221	1228	1239	rBV	141390	270034	26.98%	2.130%
31	10.700	1248	1253	1259	rVB4	3539	7607	0.76%	0.060%
32	10.824	1270	1274	1278	rBV3	1882	3254	0.33%	0.026%
33	10.924	1280	1291	1304	rBV	168818	303253	30.30%	2.392%
34	11.023	1304	1308	1309	rBV3	2101	2313	0.23%	0.018%

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Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG021321MA.M
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35	11.065	1309	1315	1326	rBV	91590	177350	17.72%	1.399%
36	11.223	1337	1342	1347	rBV3	8691	12294	1.23%	0.097%
37	11.294	1350	1354	1357	rBV3	2342	3726	0.37%	0.029%
38	12.351	1528	1534	1539	rBV5	3104	6762	0.68%	0.053%
39	12.387	1539	1540	1546	rVB	1756	1999	0.20%	0.016%
40	12.492	1555	1558	1562	rBV2	2278	2559	0.26%	0.020%
41	13.021	1646	1648	1650	rBV	1549	1736	0.17%	0.014%
42	13.074	1655	1657	1659	rBV	1666	1862	0.19%	0.015%
43	13.321	1695	1699	1702	rBV3	1864	2790	0.28%	0.022%
44	13.491	1723	1728	1733	rBV5	6292	9761	0.98%	0.077%
45	13.573	1740	1742	1745	rVB2	2002	1753	0.18%	0.014%
46	13.615	1745	1749	1754	rVB2	4053	5785	0.58%	0.046%
47	13.720	1762	1767	1770	rVB2	14707	19613	1.96%	0.155%
48	13.785	1773	1778	1782	rVB3	9408	15458	1.54%	0.122%
49	14.049	1818	1823	1827	rVB4	3213	5885	0.59%	0.046%
50	14.132	1831	1837	1842	rBV	294686	435977	43.56%	3.438%
51	14.184	1842	1846	1851	rVB	35726	53809	5.38%	0.424%
52	14.431	1881	1888	1896	rVB	317632	491272	49.09%	3.874%
53	14.602	1913	1917	1921	rBV3	2035	2762	0.28%	0.022%
54	14.731	1933	1939	1947	rBV	281378	450480	45.01%	3.553%
55	14.960	1971	1978	1986	rBV	97493	187156	18.70%	1.476%
56	15.089	1997	2000	2005	rVV5	3360	4896	0.49%	0.039%
57	15.148	2009	2010	2017	rVB6	4262	6603	0.66%	0.052%
58	15.295	2032	2035	2039	rBV3	1680	2637	0.26%	0.021%
59	15.389	2049	2051	2056	rVB3	1529	1731	0.17%	0.014%
60	15.724	2101	2108	2115	rBV	454449	650896	65.04%	5.133%
61	15.847	2123	2129	2142	rBV	146575	234175	23.40%	1.847%
62	15.965	2144	2149	2153	rVB5	6120	8857	0.88%	0.070%
63	16.400	2220	2223	2224	rBV2	2691	2641	0.26%	0.021%
64	16.499	2235	2240	2243	rBV5	3285	6493	0.65%	0.051%
65	16.664	2266	2268	2271	rBV3	1831	2360	0.24%	0.019%
66	16.693	2271	2273	2275	rVB2	3211	2982	0.30%	0.024%
67	16.922	2306	2312	2319	rBV6	13856	21701	2.17%	0.171%
68	17.357	2383	2386	2387	rBV	2374	2175	0.22%	0.017%
69	17.481	2400	2407	2417	rBV	397490	614417	61.39%	4.846%
70	17.575	2417	2423	2435	rVB2	498413	763648	76.30%	6.023%
71	17.663	2435	2438	2441	rBV3	4638	6659	0.67%	0.053%

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72	17.745	2447	2452	2457	rBV6	3858	7524	0.75%	0.059%
73	17.898	2476	2478	2482	rBV4	3495	4165	0.42%	0.033%
74	18.121	2511	2516	2523	rBV6	11613	25839	2.58%	0.204%
75	18.221	2530	2533	2534	rBV3	4069	4296	0.43%	0.034%
76	18.286	2542	2544	2545	rBV2	4452	2973	0.30%	0.023%
77	18.333	2550	2552	2553	rVV2	12029	9439	0.94%	0.074%
78	18.362	2553	2557	2563	rVV6	22415	43035	4.30%	0.339%
79	18.427	2563	2568	2571	rVV6	19377	34291	3.43%	0.270%
80	18.462	2571	2574	2578	rVB6	15181	21181	2.12%	0.167%
81	18.585	2592	2595	2599	rVB3	17498	22768	2.27%	0.180%
82	18.691	2608	2613	2618	rBV6	21395	42851	4.28%	0.338%
83	18.750	2618	2623	2630	rVV4	117589	225096	22.49%	1.775%
84	18.814	2630	2634	2641	rVB5	23476	38879	3.88%	0.307%
85	18.938	2651	2655	2664	rVB5	45762	83201	8.31%	0.656%
86	19.043	2669	2673	2674	rVV4	30027	35169	3.51%	0.277%
87	19.079	2674	2679	2685	rVB	557936	685670	68.51%	5.408%
88	19.208	2697	2701	2705	rBV3	25983	37867	3.78%	0.299%
89	19.273	2708	2712	2715	rBV6	13898	21763	2.17%	0.172%
90	19.572	2757	2763	2768	rVB	751801	983571	98.28%	7.757%
91	19.860	2806	2812	2817	rBV	667886	1000794	100.00%	7.893%
92	20.224	2871	2874	2880	rVB4	32890	41269	4.12%	0.325%
93	20.289	2880	2885	2891	rBV2	126065	180527	18.04%	1.424%
94	20.501	2918	2921	2925	rBV5	38762	46550	4.65%	0.367%
95	20.618	2938	2941	2946	rBV	40365	47138	4.71%	0.372%
96	20.812	2971	2974	2978	rVB	79668	94592	9.45%	0.746%
97	21.376	3066	3070	3075	rBV5	56798	98361	9.83%	0.776%
98	21.752	3128	3134	3141	rBV2	435985	728704	72.81%	5.747%
99	24.807	3647	3654	3663	rVB	330180	888649	88.79%	7.008%
100	25.031	3685	3692	3702	rVB2	277799	783757	78.31%	6.181%

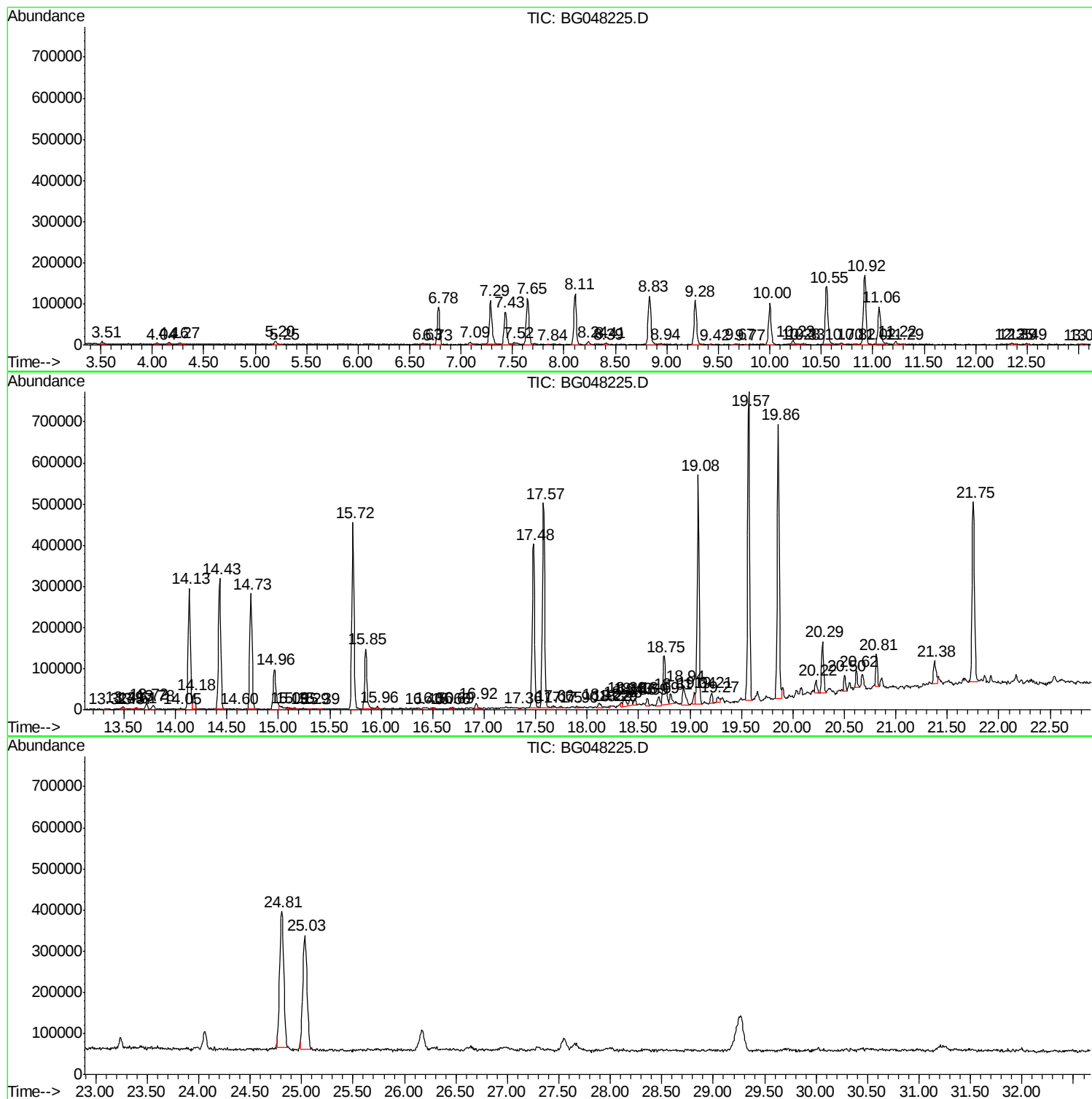
Sum of corrected areas: 12679802

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG021321MA.M
Quant Title : SVOA CALIBRATION

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



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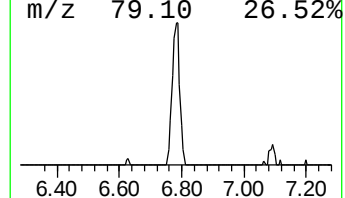
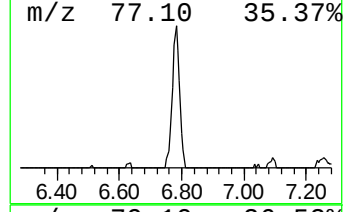
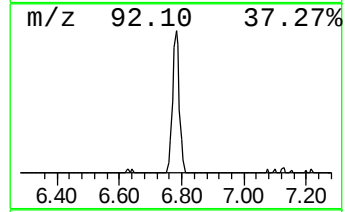
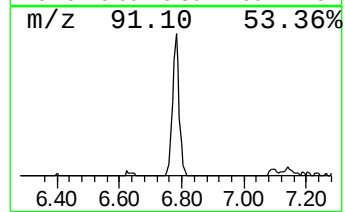
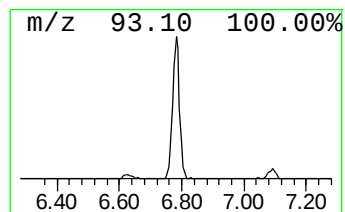
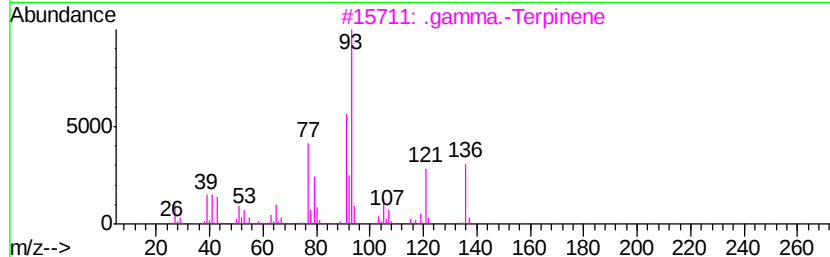
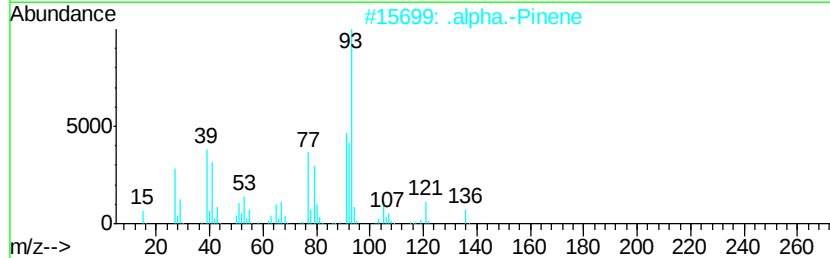
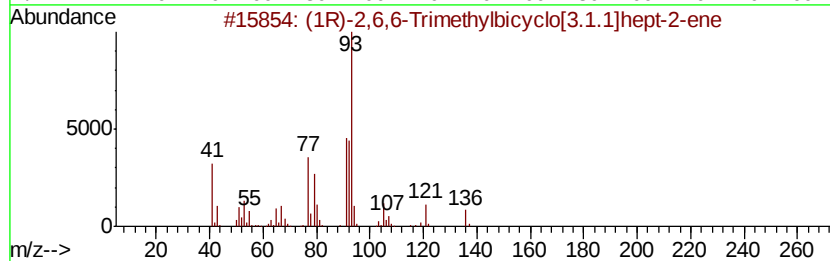
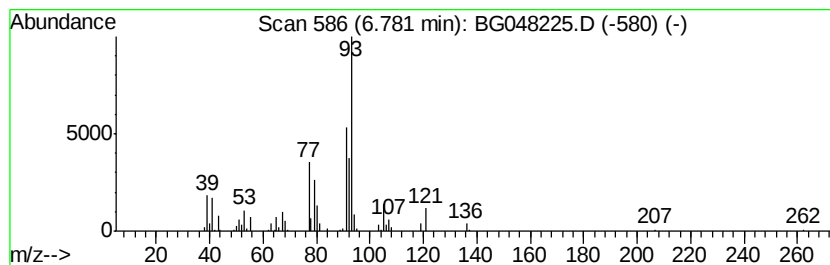
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG021321MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 (1R)-2,6,6-Trimethylbicyclo... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.78	13.32 ng/ul	143308	1,4-Dichlorobenzene-d4	8.11

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(1R)-2,6,6-Trimethylbicyclo[3.1.1]hept-2-ene	136	C10H16	007785-70-8	95
2		.alpha.-Pinene	136	C10H16	000080-56-8	95
3		.gamma.-Terpinene	136	C10H16	000099-85-4	91
4		Bicyclo[3.1.1]hept-2-ene, 3,6,6-trimethyl-	136	C10H16	004889-83-2	91
5		Tricyclo[2.2.1.0(2,6)]heptane, 1,4-dichloro-	136	C10H16	000508-32-7	91



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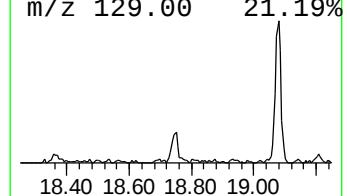
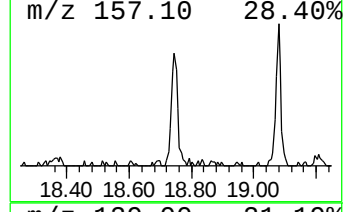
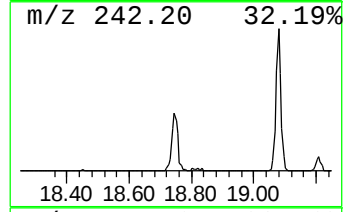
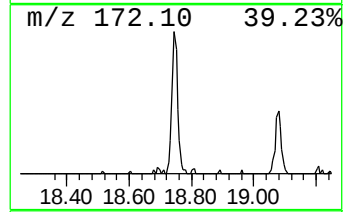
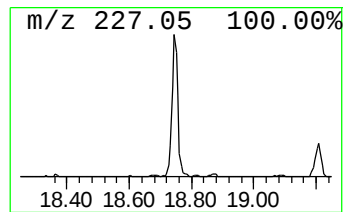
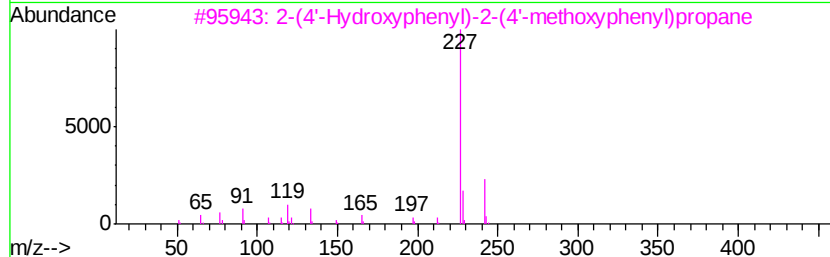
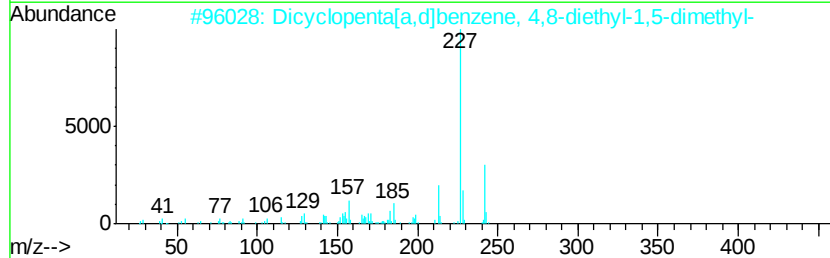
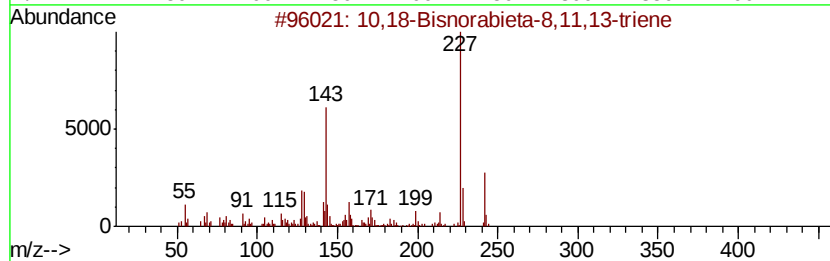
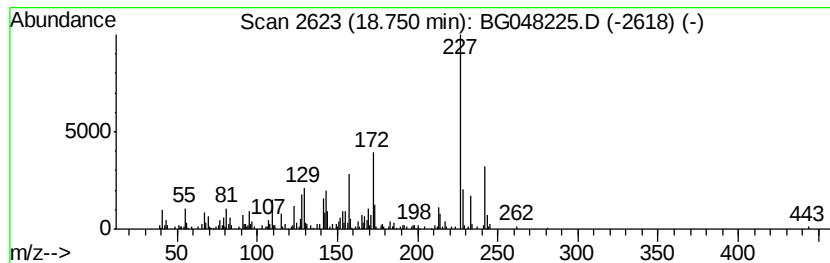
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 10,18-Bisnorabieta-8,11,13-... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.			
18.75	7.33 ng/ul	225096	Phenanthrene-d10	17.48			
Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	10,18-Bisnorabieta-8,11,13-triene	242	C18H26	032624-67-2	83		
2	Dicyclopenta[a,d]benzene, 4,8-di...	242	C18H26	1000156-41-6	70		
3	2-(4'-Hvdroxyphenyl)-2-(4'-metho...	242	C16H18O2	016530-58-8	50		
4	s-Indacene-1,7-dione, 2,3,5,6-te...	242	C16H18O2	055591-17-8	45		
5	7-Diethylamino-2-oxo-2H-chromene...	242	C14H14N2O2	332411-59-3	43		



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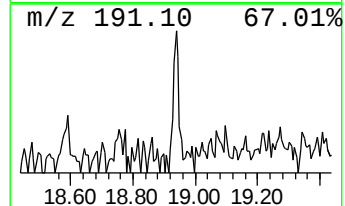
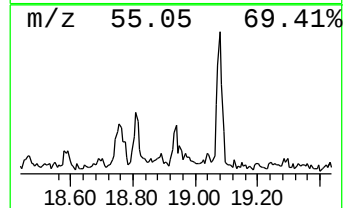
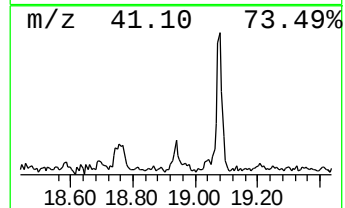
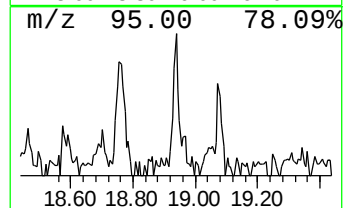
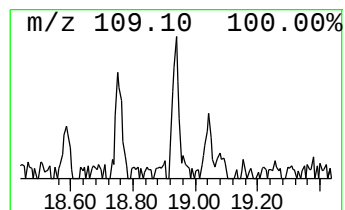
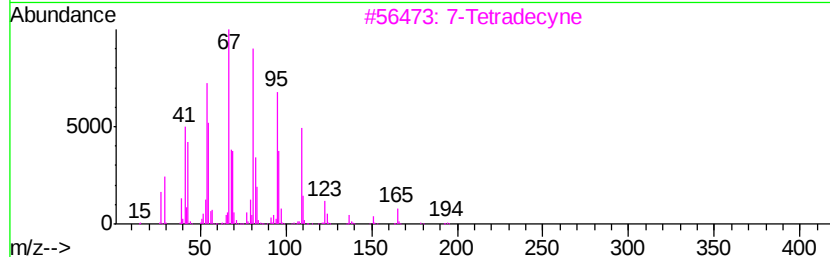
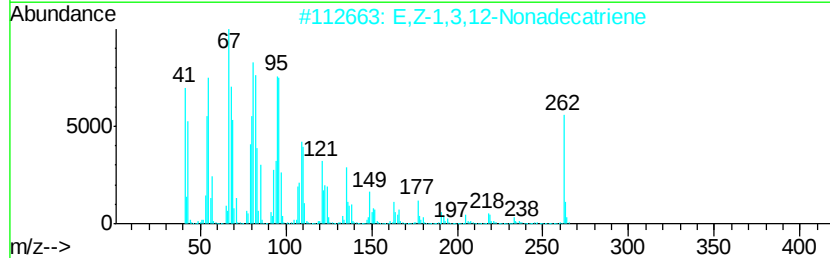
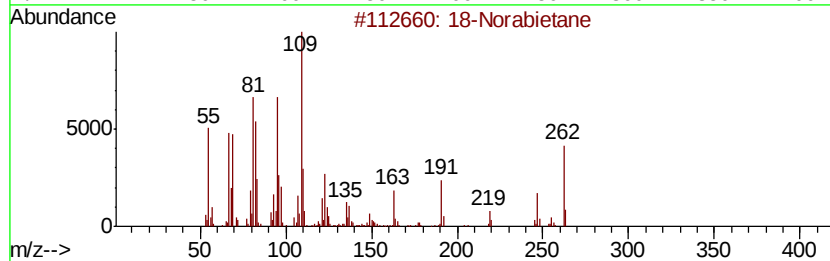
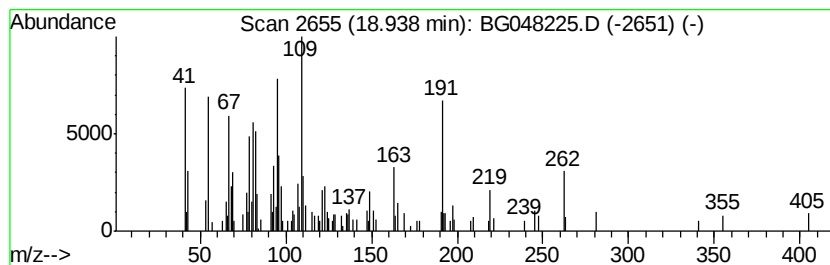
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG021321MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 18-Norabietane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.94	2.71 ng/ul	83201	Phenanthrene-d10	17.48

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	18-Norabietane		262	C19H34	1000293-16-6	55
2	E,Z-1,3,12-Nonadecatriene		262	C19H34	1000131-11-3	45
3	7-Tetradecyne		194	C14H26	035216-11-6	38
4	Cyclohexanecarboxaldehyde, 4-(hv...		142	C8H14O2	092385-32-5	27
5	Norbornane, 2-isobutyl-		152	C11H20	018127-14-5	22



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021921\
Data File : BG048225.D
Acq On : 20 Feb 2021 6:13
Operator : CG/JU
Sample : M1412-02
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBG5

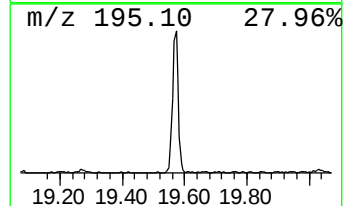
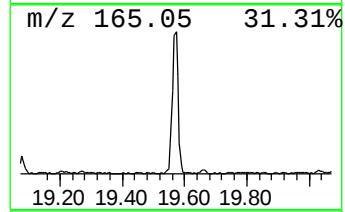
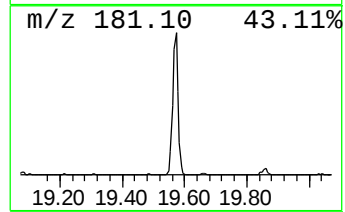
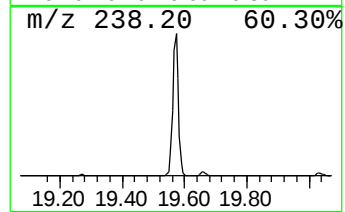
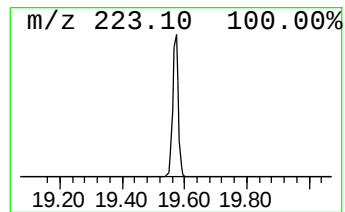
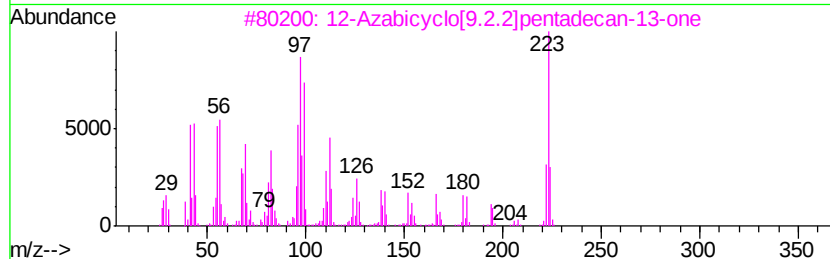
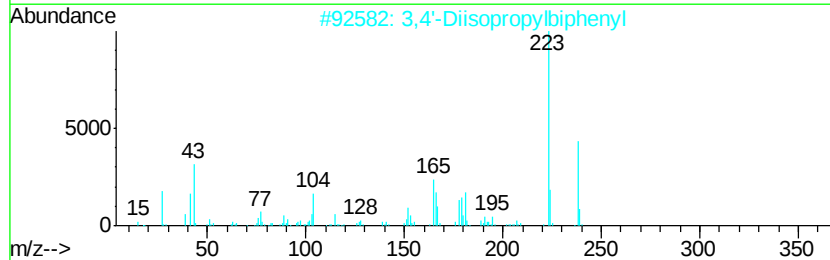
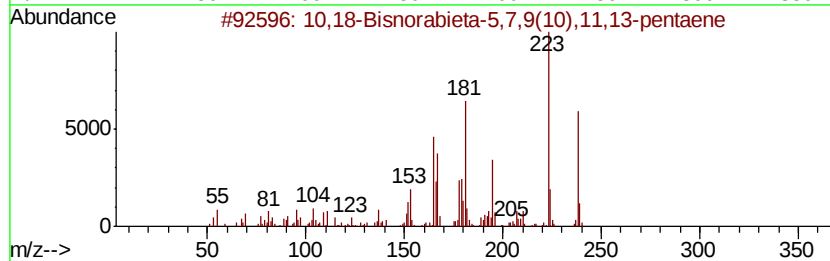
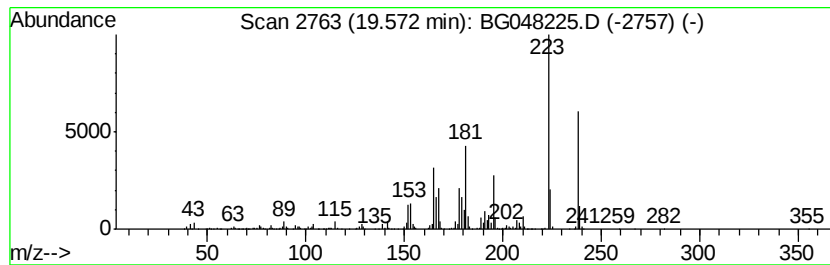
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 10,18-Bisnorabieta-5,7,9(10... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.57	32.02 ng/ul	983571	Phenanthrene-d10	17.48

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		10,18-Bisnorabieta-5,7,9(10),11,...	238	C18H22	006566-19-4	99
2		3,4'-Diisopropylbiphenyl	238	C18H22	061434-46-6	76
3		12-Azabicyclo[9.2.2]pentadecan-1...	223	C14H25NO	1000191-26-7	70
4		4,4'-Diisopropylbiphenyl	238	C18H22	018970-30-4	60
5		4,4'-Diisopropylbiphenyl	238	C18H22	018970-30-4	53



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021921\
Data File : BG048225.D
Acq On : 20 Feb 2021 6:13
Operator : CG/JU
Sample : M1412-02
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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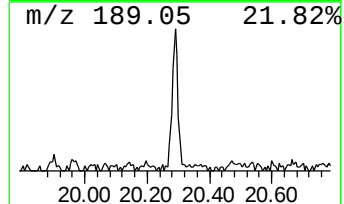
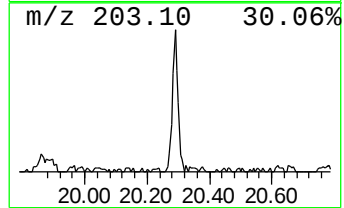
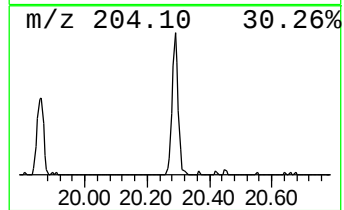
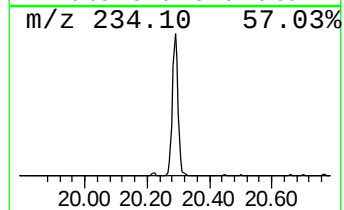
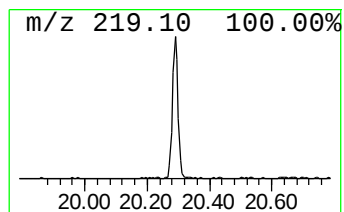
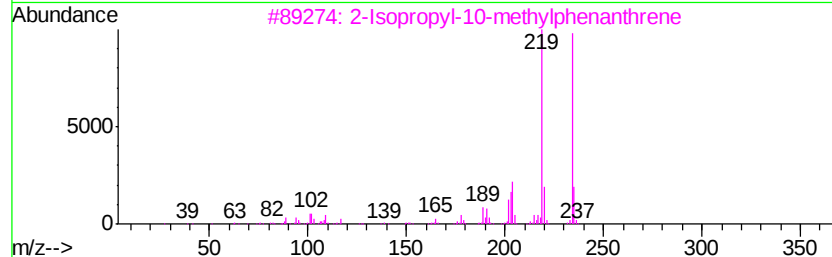
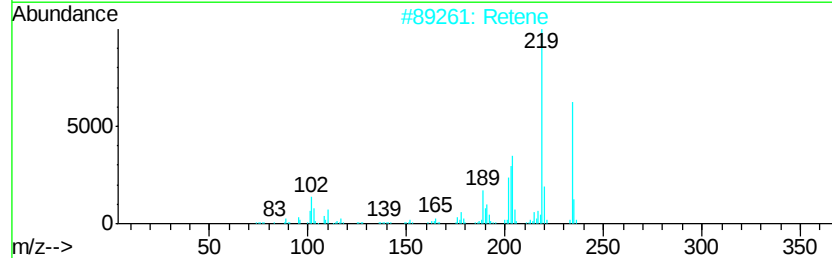
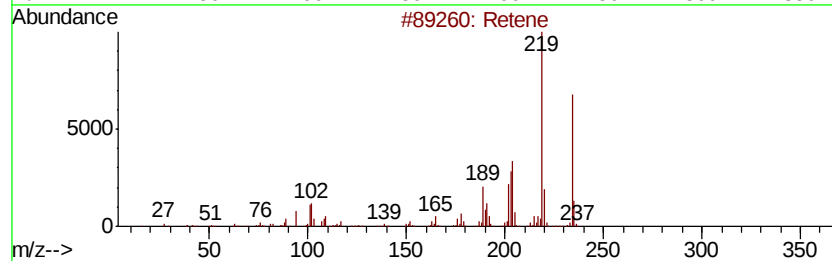
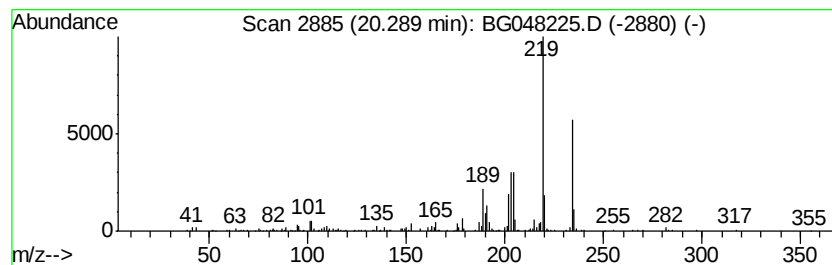
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Retene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.29	4.95 ng/ul	180527	Chrysene-d12	21.75

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Retene		234	C18H18	000483-65-8	98
2	Retene		234	C18H18	000483-65-8	95
3	2-Isopropyl-10-methylphenanthrene		234	C18H18	066552-97-4	94
4	Retene		234	C18H18	000483-65-8	93
5	Phenanthrene, 2,4,5,7-tetramethyl-		234	C18H18	007396-38-5	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021921\
Data File : BG048225.D
Acq On : 20 Feb 2021 6:13
Operator : CG/JU
Sample : M1412-02
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampled :
DBG5

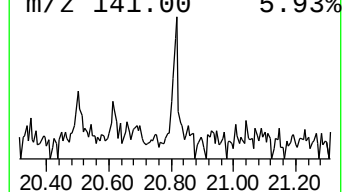
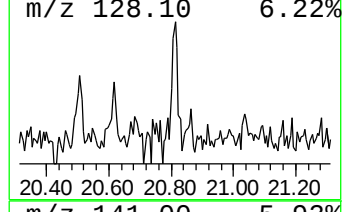
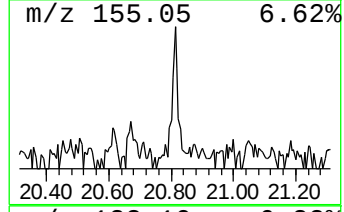
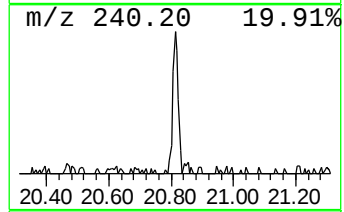
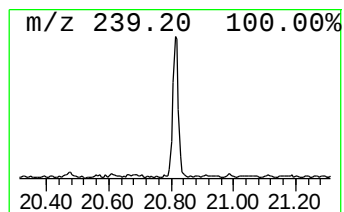
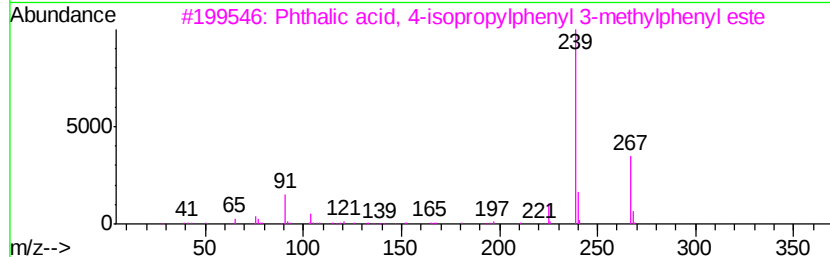
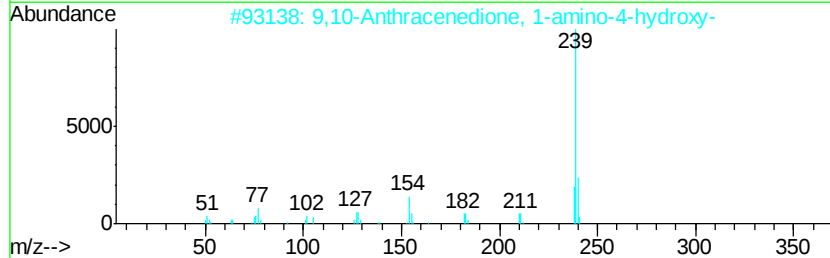
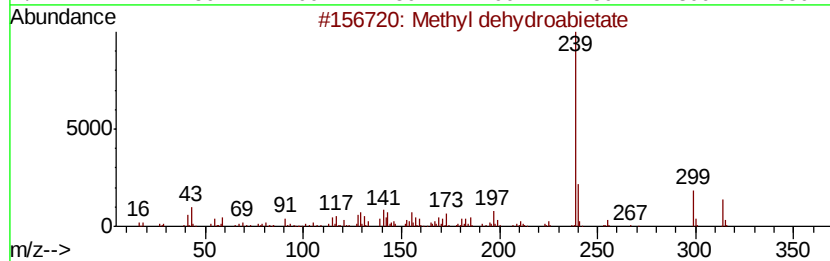
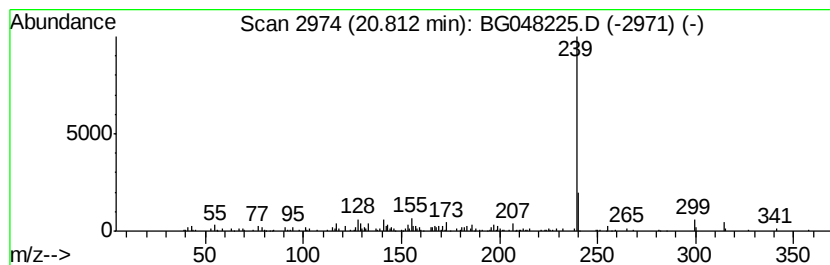
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Methyl dehydroabietate Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.81	2.60 ng/ul	94592	Chrysene-d12	21.75

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Methyl dehydroabietate	314	C21H30O2	001235-74-1	87
2		9,10-Anthracenedione, 1-amino-4-...	239	C14H9NO3	000116-85-8	64
3		Phthalic acid, 4-isopropylphenyl...	374	C24H22O4	1000315-66-1	64
4		Indole, 3-imidazo[2,1-b]thiazol...	239	C13H9N3S	292855-05-1	64
5		5H-Pyrrolo[3,4-b]pyrazine-5,7(6H...	239	C13H9N3O2	1000337-19-6	60



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021921\
Data File : BG048225.D
Acq On : 20 Feb 2021 6:13
Operator : CG/JU
Sample : M1412-02
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBGT5

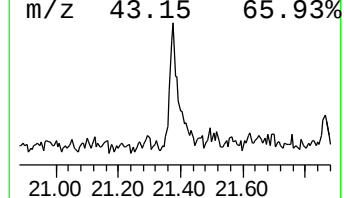
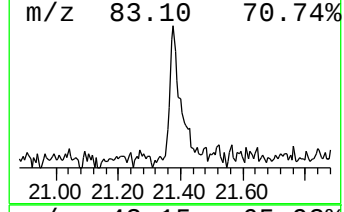
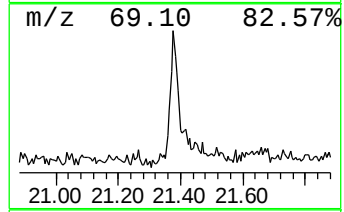
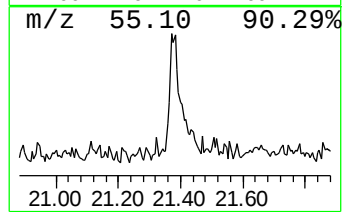
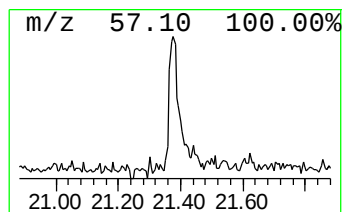
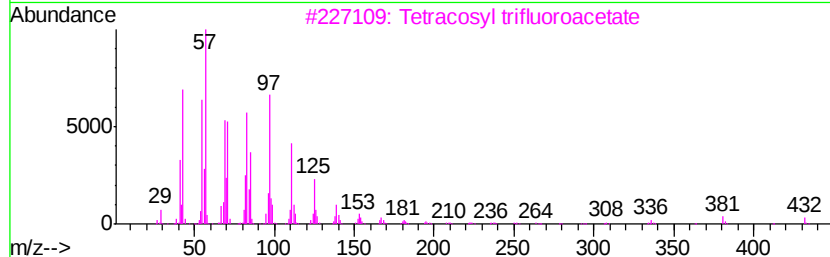
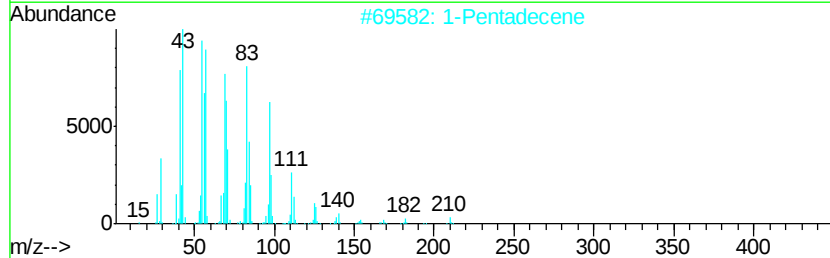
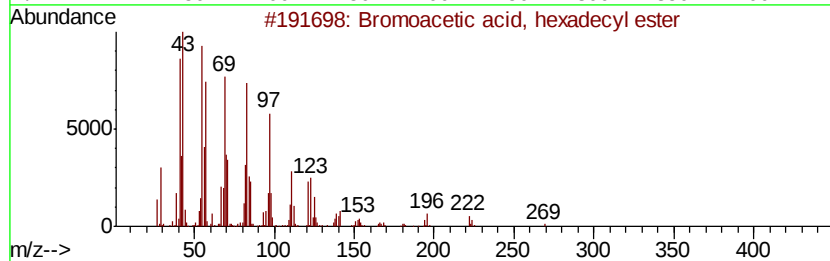
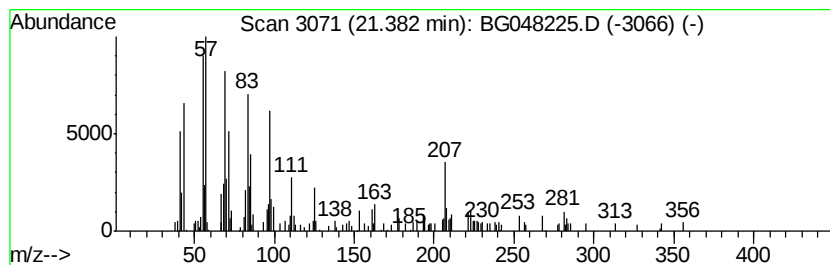
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Bromoacetic acid, hexadecyl... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.38	2.70 ng/ul	98361	Chrysene-d12	21.75

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Bromoacetic acid, hexadecyl ester	362	C18H35BrO2	005454-48-8	74
2		1-Pentadecene	210	C15H30	013360-61-7	72
3		Tetracosyl trifluoroacetate	450	C26H49F3O2	1000351-76-4	58
4		E-14-Hexadecenal	238	C16H30O	330207-53-9	56
5		Cyclohexane, 1,2-dimethyl-3-pent...	224	C16H32	062376-17-4	53



Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG021921\
Data File : BG048225.D
Acq On : 20 Feb 2021 6:13
Operator : CG/JU
Sample : M1412-02
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBGT5

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG021321MA.M
Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
(1R)-2,6,6-Trimet...	6.78	13.3	ng/ul	143308	1	8.11	215233	20.0
10,18-Bisnorabiet...	18.75	7.3	ng/ul	225096	4	17.48	614417	20.0
18-Norabietane	18.94	2.7	ng/ul	83201	4	17.48	614417	20.0
4b,8-Dimethyl-2-i...	19.08	22.3	ng/ul	685670	4	17.48	614417	20.0
10,18-Bisnorabiet...	19.57	32.0	ng/ul	983571	4	17.48	614417	20.0
Retene	20.29	5.0	ng/ul	180527	5	21.75	728704	20.0
Methyl dehydroabi...	20.81	2.6	ng/ul	94592	5	21.75	728704	20.0
Bromoacetic acid,...	21.38	2.7	ng/ul	98361	5	21.75	728704	20.0