

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111124\
 Data File : BG063342.D
 Acq On : 13 Nov 2024 2:15
 Operator : RC/JU
 Sample : P4801-10
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 GCPB1

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-BG110624.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BG063342.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.275	29	32	34	rBV4	19078	23335	0.48%	0.048%
2	3.575	80	83	87	rVB4	10689	14923	0.30%	0.031%
3	3.704	101	105	113	rBV	126534	206117	4.20%	0.422%
4	4.045	160	163	168	rVB3	9760	12509	0.25%	0.026%
5	4.280	197	203	208	rBV4	11294	16723	0.34%	0.034%
6	4.321	208	210	216	rVB5	7362	9338	0.19%	0.019%
7	4.491	234	239	243	rBV4	5515	9653	0.20%	0.020%
8	4.538	243	247	251	rBV2	60796	95934	1.96%	0.196%
9	5.002	323	326	329	rVB5	6481	7084	0.14%	0.015%
10	5.044	329	333	334	rBV4	7766	9459	0.19%	0.019%
11	5.373	385	389	391	rVB4	6412	6971	0.14%	0.014%
12	5.854	467	471	473	rBV4	6514	8921	0.18%	0.018%
13	6.712	614	617	621	rVB4	5455	7940	0.16%	0.016%
14	6.853	640	641	647	rVB3	5560	7556	0.15%	0.015%
15	7.006	661	667	674	rBV2	170859	291246	5.94%	0.596%
16	7.165	687	694	703	rBV	411366	661169	13.48%	1.354%
17	7.259	708	710	712	rVB2	8145	7145	0.15%	0.015%
18	7.364	722	728	737	rBV	483285	830936	16.94%	1.702%
19	7.429	737	739	743	rVV5	12789	16907	0.34%	0.035%
20	7.494	745	750	755	rVV9	8056	21886	0.45%	0.045%
21	7.552	757	760	764	rVB5	9671	13449	0.27%	0.028%
22	7.834	801	808	819	rBV	374988	654284	13.34%	1.340%
23	7.946	824	827	832	rVB5	15961	25271	0.52%	0.052%
24	8.093	849	852	857	rVB4	7419	11617	0.24%	0.024%
25	8.369	895	899	904	rBV5	19178	30242	0.62%	0.062%
26	8.540	920	928	941	rBV	539773	919732	18.75%	1.883%
27	8.939	992	996	997	rBV4	5040	7029	0.14%	0.014%
28	8.986	997	1004	1013	rVV	589166	1006552	20.52%	2.061%
29	9.098	1018	1023	1029	rBV5	16335	28745	0.59%	0.059%
30	9.409	1073	1076	1080	rBV5	7394	12883	0.26%	0.026%
31	9.456	1080	1084	1088	rVB5	10499	15141	0.31%	0.031%
32	9.515	1090	1094	1100	rBV4	24062	35561	0.72%	0.073%
33	9.585	1102	1106	1108	rBV3	8570	13448	0.27%	0.028%
34	9.709	1119	1127	1136	rBV	528063	951938	19.40%	1.949%
35	9.867	1152	1154	1158	rVB2	6329	8268	0.17%	0.017%
36	10.050	1175	1185	1192	rBV5	94236	250881	5.11%	0.514%

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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

Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110624.MA.M
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37	10.120	1192	1197	1202	rBV2	159775	277611	5.66%	0.568%
38	10.249	1211	1219	1230	rBV	823023	1487705	30.32%	3.046%
39	10.402	1241	1245	1246	rBV4	11600	13200	0.27%	0.027%
40	10.619	1275	1282	1289	rBV	604562	1042643	21.25%	2.135%
41	10.755	1298	1305	1319	rBV2	545717	1154306	23.53%	2.364%
42	11.136	1368	1370	1374	rVB6	10749	13445	0.27%	0.028%
43	11.442	1416	1422	1425	rBV6	10191	19073	0.39%	0.039%
44	11.665	1456	1460	1462	rBV3	15969	17803	0.36%	0.036%
45	11.812	1480	1485	1492	rBV3	30614	57814	1.18%	0.118%
46	11.900	1496	1500	1505	rBV5	10330	19276	0.39%	0.039%
47	11.941	1505	1507	1512	rVB4	5332	8313	0.17%	0.017%
48	12.006	1515	1518	1524	rBV5	8201	12054	0.25%	0.025%
49	12.124	1534	1538	1540	rBV3	8989	11035	0.22%	0.023%
50	12.176	1542	1547	1551	rBV2	52730	91065	1.86%	0.186%
51	12.288	1561	1566	1572	rBV4	35893	71170	1.45%	0.146%
52	12.417	1584	1588	1590	rBV5	4369	7069	0.14%	0.014%
53	12.505	1593	1603	1612	rVB3	50296	131029	2.67%	0.268%
54	12.588	1612	1617	1618	rBV3	5954	8558	0.17%	0.018%
55	12.676	1628	1632	1641	rVB3	24124	43631	0.89%	0.089%
56	12.752	1641	1645	1648	rBV4	6393	8120	0.17%	0.017%
57	12.817	1652	1656	1659	rVB5	9775	12815	0.26%	0.026%
58	12.870	1659	1665	1669	rBV4	45991	78390	1.60%	0.161%
59	13.070	1697	1699	1703	rVB5	9934	11725	0.24%	0.024%
60	13.305	1735	1739	1744	rBV4	22523	34940	0.71%	0.072%
61	13.363	1744	1749	1756	rVV	270705	415447	8.47%	0.851%
62	13.434	1756	1761	1764	rBV5	11126	20797	0.42%	0.043%
63	13.534	1773	1778	1782	rVB8	16793	32678	0.67%	0.067%
64	13.604	1786	1790	1792	rBV4	6853	8027	0.16%	0.016%
65	13.651	1792	1798	1804	rBV9	29887	62514	1.27%	0.128%
66	13.798	1817	1823	1829	rVB6	26212	42770	0.87%	0.088%
67	13.874	1830	1836	1849	rVB	1666183	2445070	49.84%	5.007%
68	14.115	1873	1877	1878	rBV3	18263	19286	0.39%	0.039%
69	14.162	1878	1885	1898	rVV2	1823955	3578720	72.94%	7.328%
70	14.280	1901	1905	1913	rVB7	19437	39143	0.80%	0.080%
71	14.409	1922	1927	1931	rBV7	27990	54243	1.11%	0.111%
72	14.468	1931	1937	1944	rBV	1092532	1710441	34.86%	3.503%
73	14.685	1968	1974	1982	rBV3	157689	305729	6.23%	0.626%
74	14.832	1997	1999	2002	rVB4	11386	10298	0.21%	0.021%
75	15.085	2037	2042	2043	rBV5	11901	19108	0.39%	0.039%
76	15.337	2081	2085	2089	rBV4	19467	37172	0.76%	0.076%

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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	15.414	2093	2098	2099	rBV4	11492	15725	0.32%	0.032%
78	15.467	2100	2107	2113	rBV	2542417	3649076	74.38%	7.472%
79	15.590	2122	2128	2134	rBV	8433315	1182091	24.09%	2.421%
80	15.825	2166	2168	2174	rVB4	26125	34931	0.71%	0.072%
81	16.471	2277	2278	2281	rVB3	9210	7024	0.14%	0.014%
82	16.759	2325	2327	2330	rVB4	14326	9675	0.20%	0.020%
83	17.141	2388	2392	2396	rVV2	138191	162714	3.32%	0.333%
84	17.223	2398	2406	2413	rBV	1444657	2180429	44.44%	4.465%
85	17.323	2416	2423	2431	rVB	2690687	4083077	83.22%	8.361%
86	17.500	2450	2453	2456	rBV5	14166	24219	0.49%	0.050%
87	17.629	2473	2475	2479	rVB3	19218	16923	0.34%	0.035%
88	17.717	2485	2490	2493	rBV	48976	79158	1.61%	0.162%
89	17.893	2516	2520	2525	rVB7	77639	114726	2.34%	0.235%
90	18.122	2555	2559	2567	rBV5	88916	146932	2.99%	0.301%
91	18.340	2591	2596	2612	rBV	1208434	2169065	44.21%	4.442%
92	18.545	2628	2631	2636	rVB2	105484	133830	2.73%	0.274%
93	19.068	2717	2720	2726	rVB7	53277	64475	1.31%	0.132%
94	19.174	2736	2738	2747	rVB5	50384	87058	1.77%	0.178%
95	19.250	2748	2751	2754	rBV3	34674	54500	1.11%	0.112%
96	19.403	2774	2777	2782	rVB7	38765	52851	1.08%	0.108%
97	19.621	2808	2814	2821	rVB	3498964	4850070	98.86%	9.932%
98	21.471	3123	3129	3138	rBV	1553769	2485108	50.65%	5.089%
99	24.303	3601	3611	3623	rVB	1894554	4906155	100.00%	10.046%
100	24.509	3637	3646	3658	rVB	1026234	2639722	53.80%	5.405%

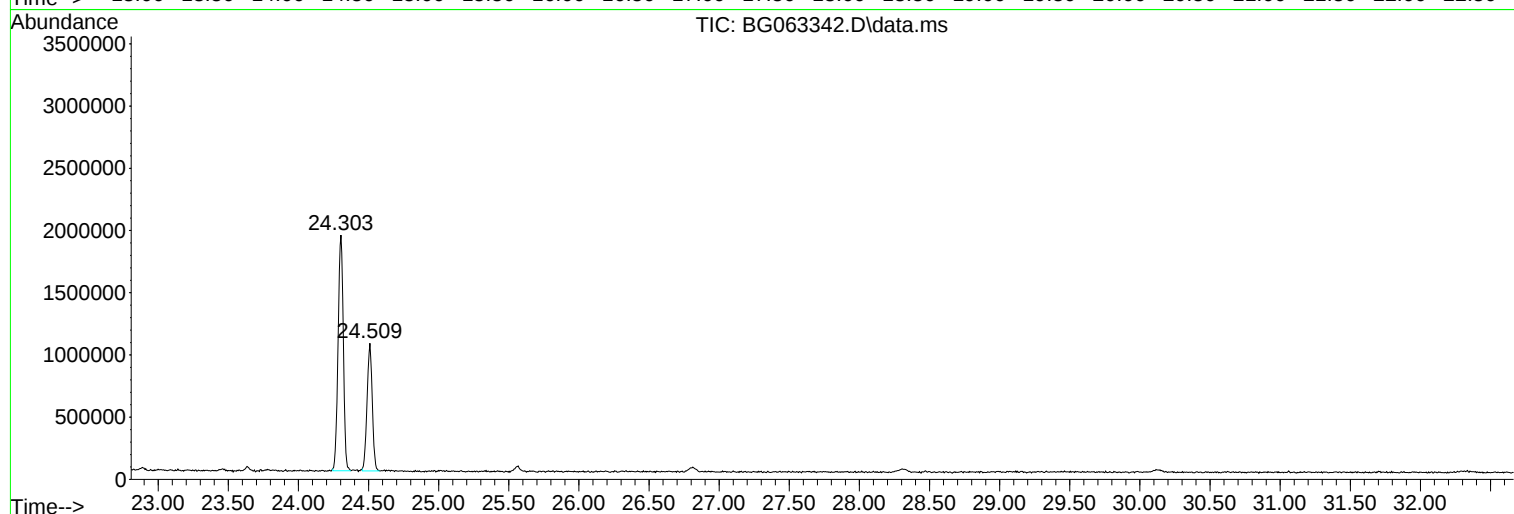
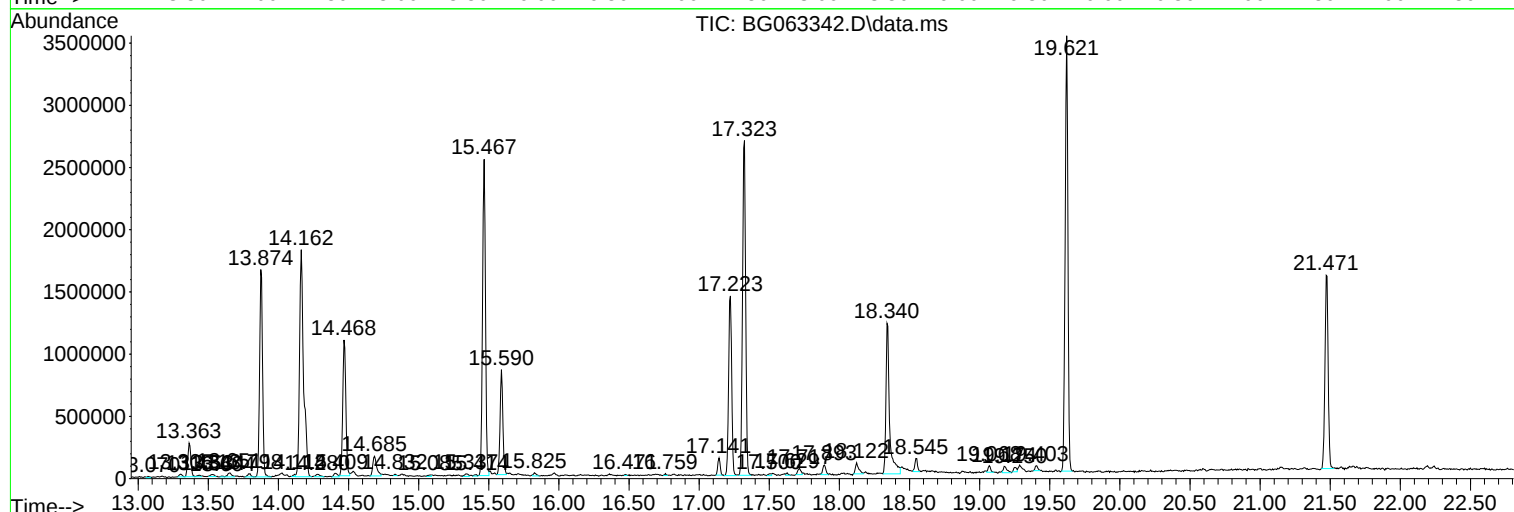
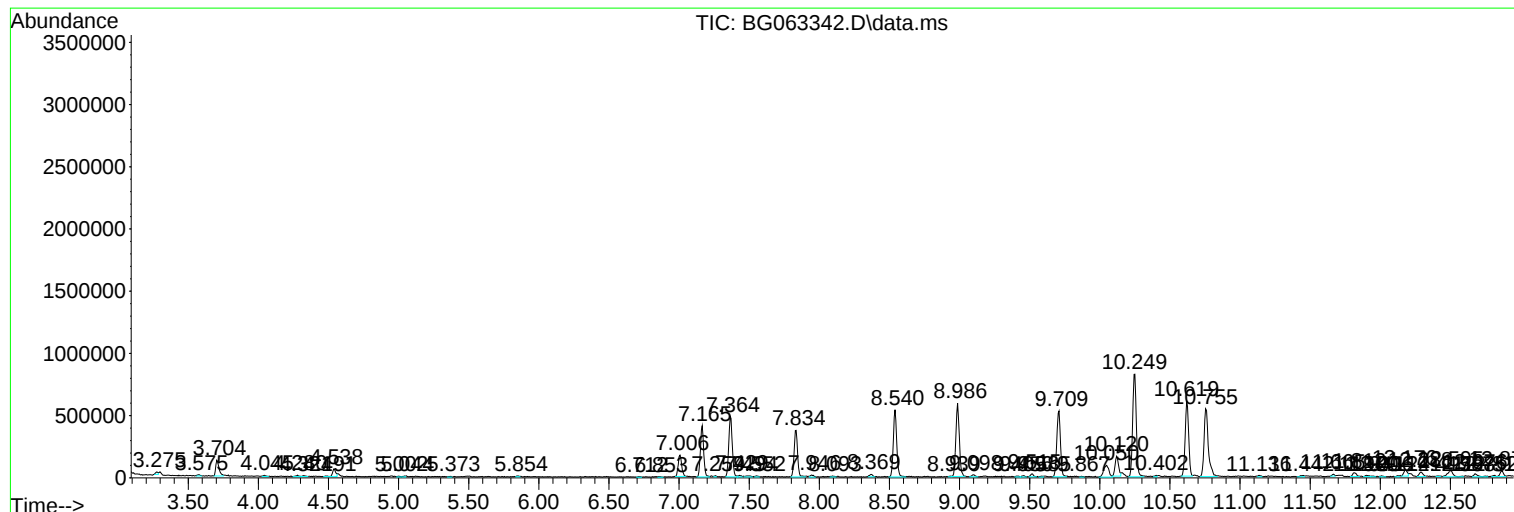
Sum of corrected areas: 48834560

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ALS Vial : 22 Sample Multiplier: 1

Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110624.MA.M
Quant Title : SVOA CALIBRATION

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



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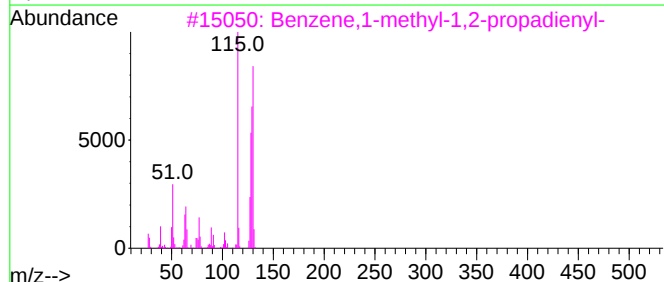
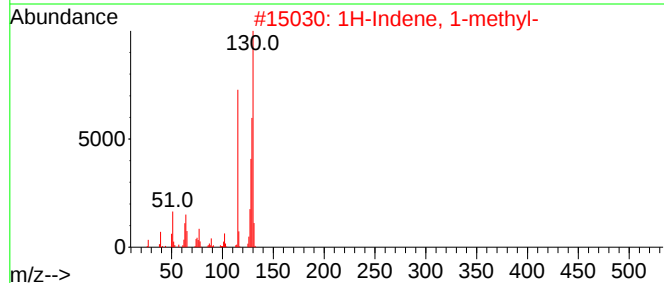
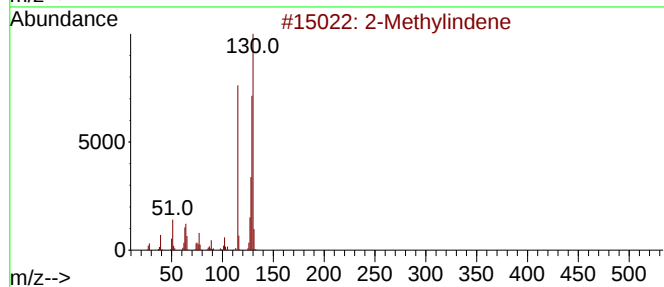
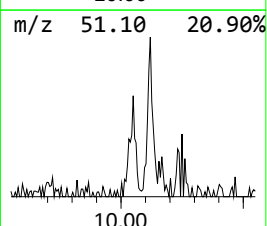
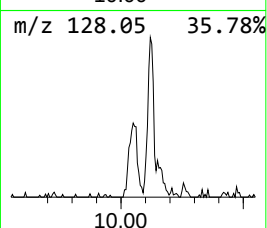
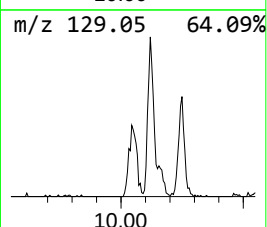
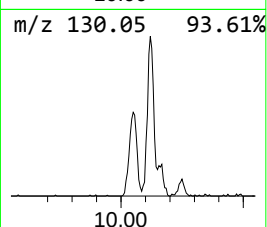
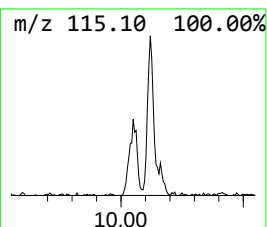
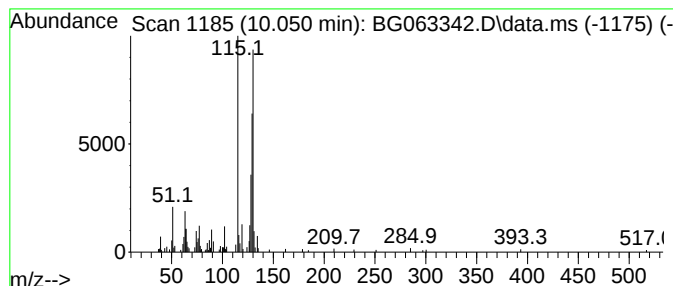
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110624.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 2-Methylindene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.050	4.81 ng/ul	250881	Naphthalene-d8	10.619

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Methylindene	130	C10H10	002177-47-1	93
2			1H-Indene, 1-methyl-	130	C10H10	000767-59-9	93
3			Benzene,1-methyl-1,2-propadienyl-	130	C10H10	022433-39-2	91
4			2-Methylindene	130	C10H10	002177-47-1	91
5			1H-Indene, 1-methyl-	130	C10H10	000767-59-9	90



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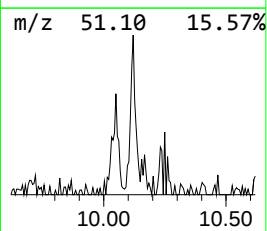
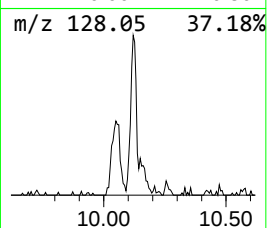
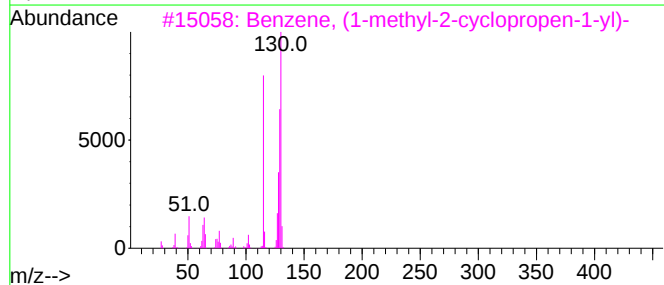
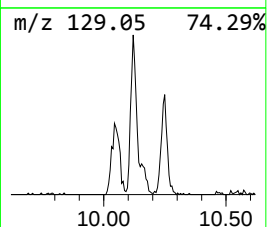
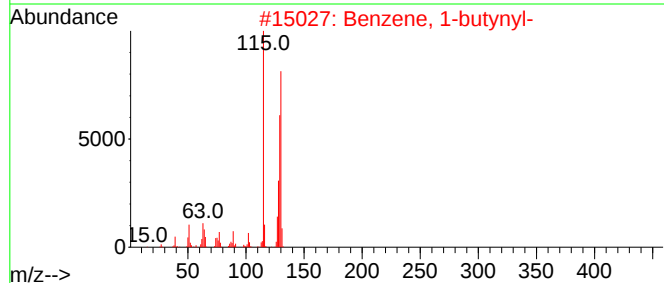
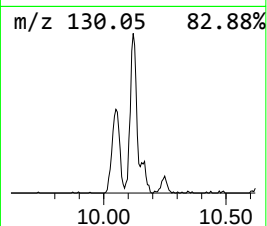
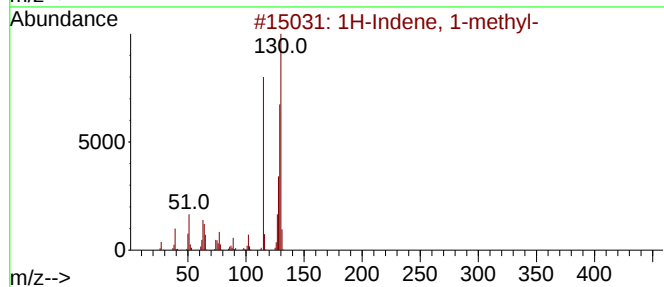
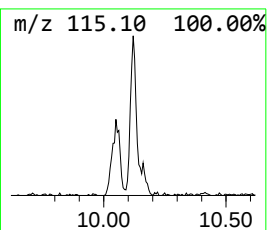
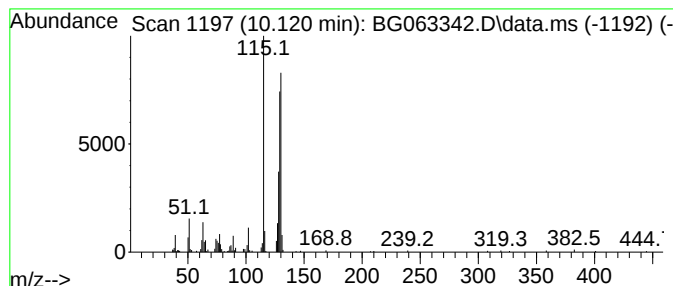
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110624.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 1H-Indene, 1-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.120	5.33 ng/ul	277611	Naphthalene-d8	10.619

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 1-methyl-	130	C10H10	000767-59-9	93
2			Benzene, 1-butyryl-	130	C10H10	000622-76-4	93
3			Benzene, (1-methyl-2-cyclopropen-1-yl)-	130	C10H10	065051-83-4	93
4			2-Methylindene	130	C10H10	002177-47-1	90
5			2-Methylindene	130	C10H10	002177-47-1	89



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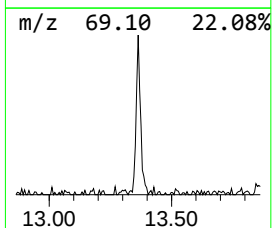
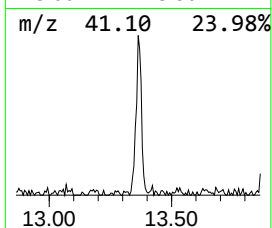
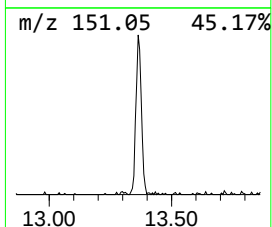
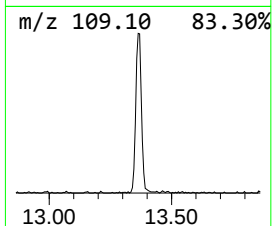
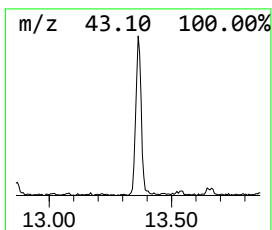
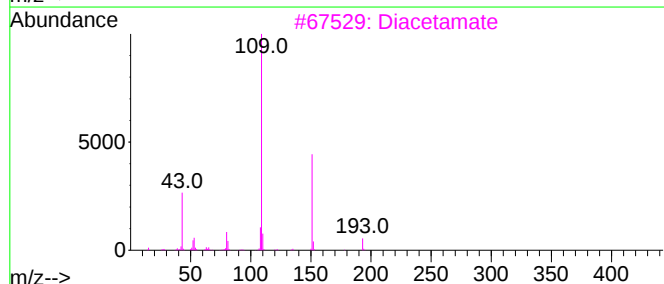
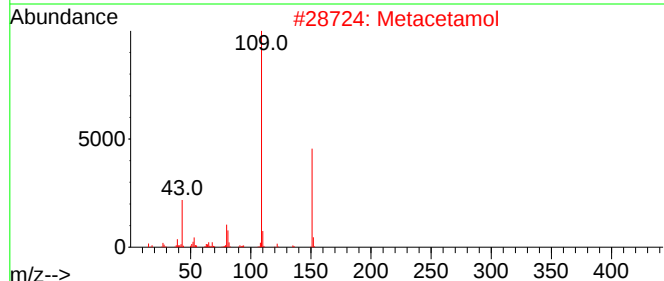
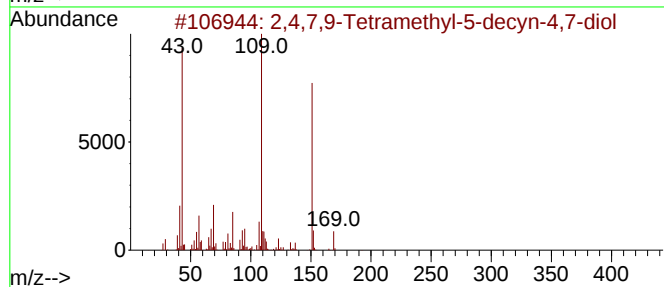
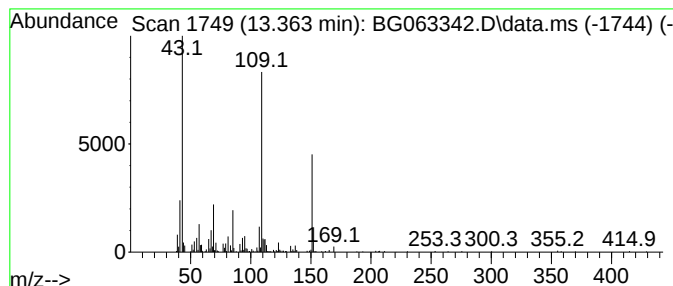
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110624.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 5 2,4,7,9-Tetramethyl-5-decyn... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.363	4.86 ng/ul	415447	Acenaphthene-d10	14.468

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4,7,9-Tetramethyl-5-decyn-4,7-...	226	C14H26O2	000126-86-3	87	
2	Metacetamol	151	C8H9NO2	000621-42-1	58	
3	Diacetamate	193	C10H11NO3	002623-33-8	53	
4	Acetaminophen	151	C8H9NO2	000103-90-2	53	
5	Metacetamol	151	C8H9NO2	000621-42-1	53	



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111124\
Data File : BG063342.D
Acq On : 13 Nov 2024 2:15
Operator : RC/JU
Sample : P4801-10
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
GCPB1

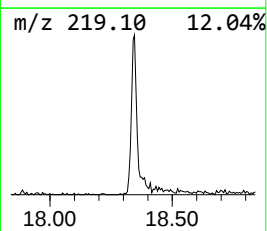
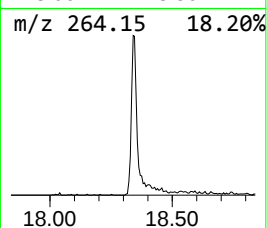
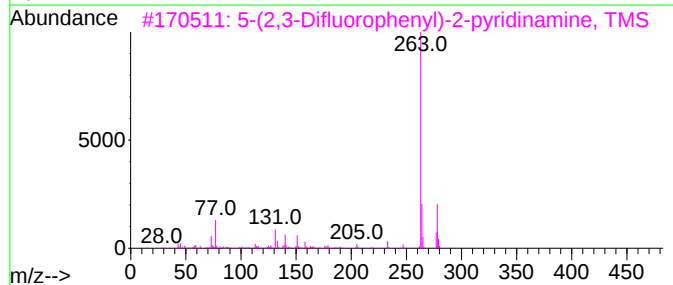
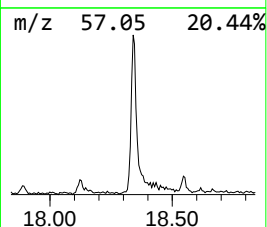
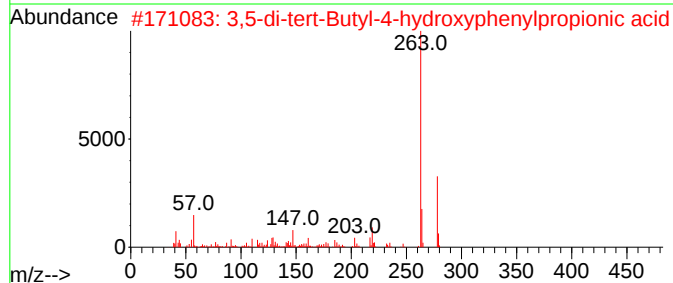
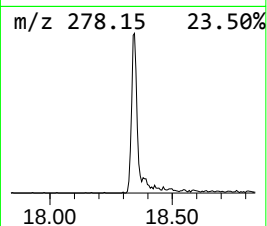
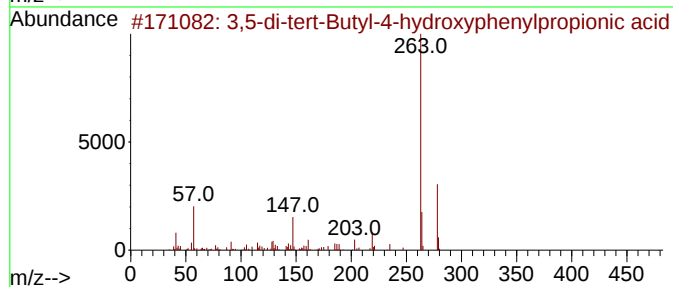
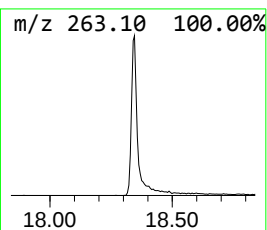
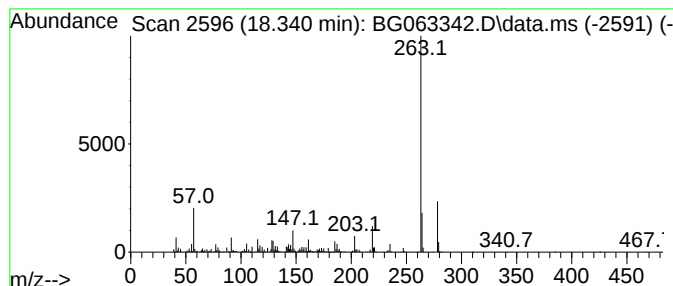
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 3,5-di-tert-Butyl-4-hydroxy... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.340	19.90 ng/ul	2169070	Phenanthrene-d10	17.224

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3,5-di-tert-Butyl-4-hydroxypheny...	278	C17H26O3	020170-32-5	94
2			3,5-di-tert-Butyl-4-hydroxypheny...	278	C17H26O3	020170-32-5	94
3			5-(2,3-Difluorophenyl)-2-pyridin...	278	C14H16F2N2Si	1000504-49-4	72
4			Cyanic acid, (1-methylethylidene...	278	C17H14N2O2	001156-51-0	64
5			6,7-Dimethoxy-4(1H)-quinazolinon...	278	C13H18N2O3Si	1000479-65-5	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111124\
Data File : BG063342.D
Acq On : 13 Nov 2024 2:15
Operator : RC/JU
Sample : P4801-10
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
GCPB1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110624.MA.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
2-Methylindene	10.050	4.8	ng/ul	250881	2	10.619	1042640	20.0
1H-Indene, 1-me...	10.120	5.3	ng/ul	277611	2	10.619	1042640	20.0
2,4,7,9-Tetrame...	13.363	4.9	ng/ul	415447	3	14.468	1710440	20.0
3,5-di-tert-But...	18.340	19.9	ng/ul	2169070	4	17.224	2180430	20.0