

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021921\
 Data File : BG048254.D
 Acq On : 21 Feb 2021 3:33
 Operator : CG/JU
 Sample : M1415-13
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 DBH27

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	3.515	27	30	34	rVB3	4338	5587	0.74%	0.054%
2	3.809	75	80	84	rBV2	3413	6069	0.80%	0.059%
3	4.050	117	121	125	rBV6	3177	5684	0.75%	0.055%
4	4.167	138	141	146	rVB2	3298	4285	0.56%	0.041%
5	4.267	154	158	165	rVV	3852	5544	0.73%	0.054%
6	4.391	178	179	184	rBV3	1091	1522	0.20%	0.015%
7	4.649	220	223	225	rBV2	1240	1412	0.19%	0.014%
8	4.678	225	228	232	rVB2	1283	1687	0.22%	0.016%
9	4.972	275	278	281	rBV2	1144	1559	0.21%	0.015%
10	5.190	312	315	319	rBV5	3000	4568	0.60%	0.044%
11	5.472	358	363	366	rBV2	1032	1555	0.20%	0.015%
12	5.624	388	389	392	rBV3	1691	1571	0.21%	0.015%
13	5.918	436	439	442	rBV	1027	1510	0.20%	0.015%
14	6.177	481	483	489	rVB2	1791	2535	0.33%	0.025%
15	6.294	499	503	504	rBV	1465	1517	0.20%	0.015%
16	6.776	583	585	590	rVB2	3501	4951	0.65%	0.048%
17	7.093	635	639	642	rVB	2620	3276	0.43%	0.032%
18	7.258	662	667	670	rBV3	6430	9972	1.31%	0.097%
19	7.316	670	677	690	rVV2	124744	254582	33.51%	2.466%
20	7.440	691	698	706	rBV	81693	136400	17.95%	1.321%
21	7.522	706	712	718	rVB3	5040	8701	1.15%	0.084%
22	7.604	723	726	727	rBV	1541	1570	0.21%	0.015%
23	7.657	730	735	740	rBV	114409	205748	27.08%	1.993%
24	8.116	806	813	819	rBV	159712	271971	35.80%	2.634%
25	8.239	831	834	840	rVB4	4163	6061	0.80%	0.059%
26	8.386	856	859	861	rBV	2338	3013	0.40%	0.029%
27	8.486	873	876	879	rBV2	1918	1832	0.24%	0.018%
28	8.856	931	939	949	rBV3	141364	249309	32.82%	2.414%
29	8.944	949	954	959	rBV4	3478	5173	0.68%	0.050%
30	9.285	1004	1012	1021	rBV	110018	202295	26.63%	1.959%
31	9.349	1021	1023	1024	rBV	2452	1873	0.25%	0.018%
32	9.655	1072	1075	1079	rVB	2602	2781	0.37%	0.027%
33	9.843	1102	1107	1110	rBV4	3784	5230	0.69%	0.051%
34	10.013	1127	1136	1144	rBV	104559	196306	25.84%	1.901%

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35	10.278	1174	1181	1186	rBV6	4164	10012	1.32%	0.097%
36	10.336	1188	1191	1196	rVV5	9416	13824	1.82%	0.134%
37	10.389	1196	1200	1203	rVB2	1438	2458	0.32%	0.024%
38	10.572	1224	1231	1239	rBV	153595	280830	36.97%	2.720%
39	10.707	1247	1254	1259	rBV2	11490	22470	2.96%	0.218%
40	10.830	1271	1275	1276	rBV2	1637	1699	0.22%	0.016%
41	10.848	1276	1278	1284	rVB4	1338	2445	0.32%	0.024%
42	10.930	1285	1292	1299	rBV	213028	380334	50.06%	3.683%
43	10.983	1299	1301	1307	rVB3	6319	8037	1.06%	0.078%
44	11.083	1311	1318	1325	rBV	66344	129203	17.01%	1.251%
45	11.394	1368	1371	1372	rVB3	1968	1496	0.20%	0.014%
46	11.476	1378	1385	1390	rVB4	4223	7943	1.05%	0.077%
47	11.564	1396	1400	1401	rBV2	1559	1964	0.26%	0.019%
48	11.670	1414	1418	1421	rBV	1648	1810	0.24%	0.018%
49	11.917	1458	1460	1465	rVB2	1245	1879	0.25%	0.018%
50	12.217	1508	1511	1514	rBV	953	1463	0.19%	0.014%
51	12.505	1557	1560	1563	rBV	3483	4302	0.57%	0.042%
52	12.651	1581	1585	1587	rBV2	3716	4620	0.61%	0.045%
53	12.704	1592	1594	1599	rVB2	1649	1611	0.21%	0.016%
54	12.804	1607	1611	1614	rVB3	1273	1636	0.22%	0.016%
55	12.928	1628	1632	1635	rBV2	1141	1745	0.23%	0.017%
56	13.098	1659	1661	1665	rVB4	1822	2145	0.28%	0.021%
57	13.245	1684	1686	1689	rVB	1471	1438	0.19%	0.014%
58	13.345	1700	1703	1705	rVB3	2728	2253	0.30%	0.022%
59	13.398	1708	1712	1714	rBV	1108	1516	0.20%	0.015%
60	13.497	1728	1729	1734	rBV3	1455	2163	0.28%	0.021%
61	13.544	1734	1737	1738	rBV2	1562	1820	0.24%	0.018%
62	13.633	1745	1752	1756	rBV	83978	137741	18.13%	1.334%
63	13.674	1756	1759	1766	rVV2	24687	40996	5.40%	0.397%
64	13.727	1766	1768	1772	rVB3	1909	2185	0.29%	0.021%
65	13.809	1781	1782	1785	rBV3	2262	2198	0.29%	0.021%
66	13.844	1786	1788	1791	rVB3	1602	1497	0.20%	0.014%
67	13.868	1791	1792	1795	rBV3	1298	1491	0.20%	0.014%
68	13.921	1798	1801	1803	rBV3	1057	1541	0.20%	0.015%
69	14.103	1829	1832	1833	rBV	1656	1626	0.21%	0.016%
70	14.144	1833	1839	1844	rVV	268037	400907	52.77%	3.883%
71	14.191	1844	1847	1858	rVB	79890	112182	14.77%	1.086%

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72	14.273	1858	1861	1863	rBV4	1851	2022	0.27%	0.020%
73	14.367	1873	1877	1882	rVB4	3057	4585	0.60%	0.044%
74	14.438	1882	1889	1896	rBV	301494	465404	61.26%	4.507%
75	14.549	1905	1908	1909	rVB3	1667	1657	0.22%	0.016%
76	14.567	1909	1911	1914	rBV4	2584	3076	0.40%	0.030%
77	14.608	1914	1918	1925	rVB4	7251	11778	1.55%	0.114%
78	14.743	1934	1941	1948	rBV	359572	556854	73.30%	5.393%
79	14.831	1954	1956	1958	rBV3	1487	1433	0.19%	0.014%
80	14.919	1970	1971	1975	rVB2	1694	1672	0.22%	0.016%
81	15.007	1979	1986	2001	rBV	106633	214138	28.19%	2.074%
82	15.119	2001	2005	2010	rBV7	9610	15745	2.07%	0.152%
83	15.342	2041	2043	2045	rBV2	2637	2571	0.34%	0.025%
84	15.413	2053	2055	2057	rBV3	2290	2637	0.35%	0.026%
85	15.548	2075	2078	2083	rVB4	7980	9079	1.20%	0.088%
86	15.736	2104	2110	2117	rBV	367586	573335	75.47%	5.553%
87	15.865	2125	2132	2140	rBV	131044	206673	27.20%	2.002%
88	16.388	2220	2221	2223	rBV2	4333	2970	0.39%	0.029%
89	16.623	2259	2261	2263	rBV3	4145	2646	0.35%	0.026%
90	17.493	2402	2409	2415	rBV	416490	658521	86.68%	6.378%
91	17.587	2420	2425	2432	rVB	397818	597978	78.71%	5.791%
92	18.592	2592	2596	2600	rVB2	76904	102653	13.51%	0.994%
93	18.762	2620	2625	2630	rBV	350204	488170	64.26%	4.728%
94	18.944	2651	2656	2664	rBV	347379	498885	65.67%	4.832%
95	19.085	2676	2680	2685	rVB	255367	314489	41.40%	3.046%
96	19.573	2760	2763	2768	rVB	95928	127092	16.73%	1.231%
97	19.872	2809	2814	2821	rVB	453978	686150	90.32%	6.645%
98	20.301	2883	2887	2892	rVB	375410	475189	62.55%	4.602%
99	20.818	2972	2975	2980	rVB	256728	325340	42.82%	3.151%
100	21.770	3132	3137	3144	rVB2	444496	759706	100.00%	7.358%

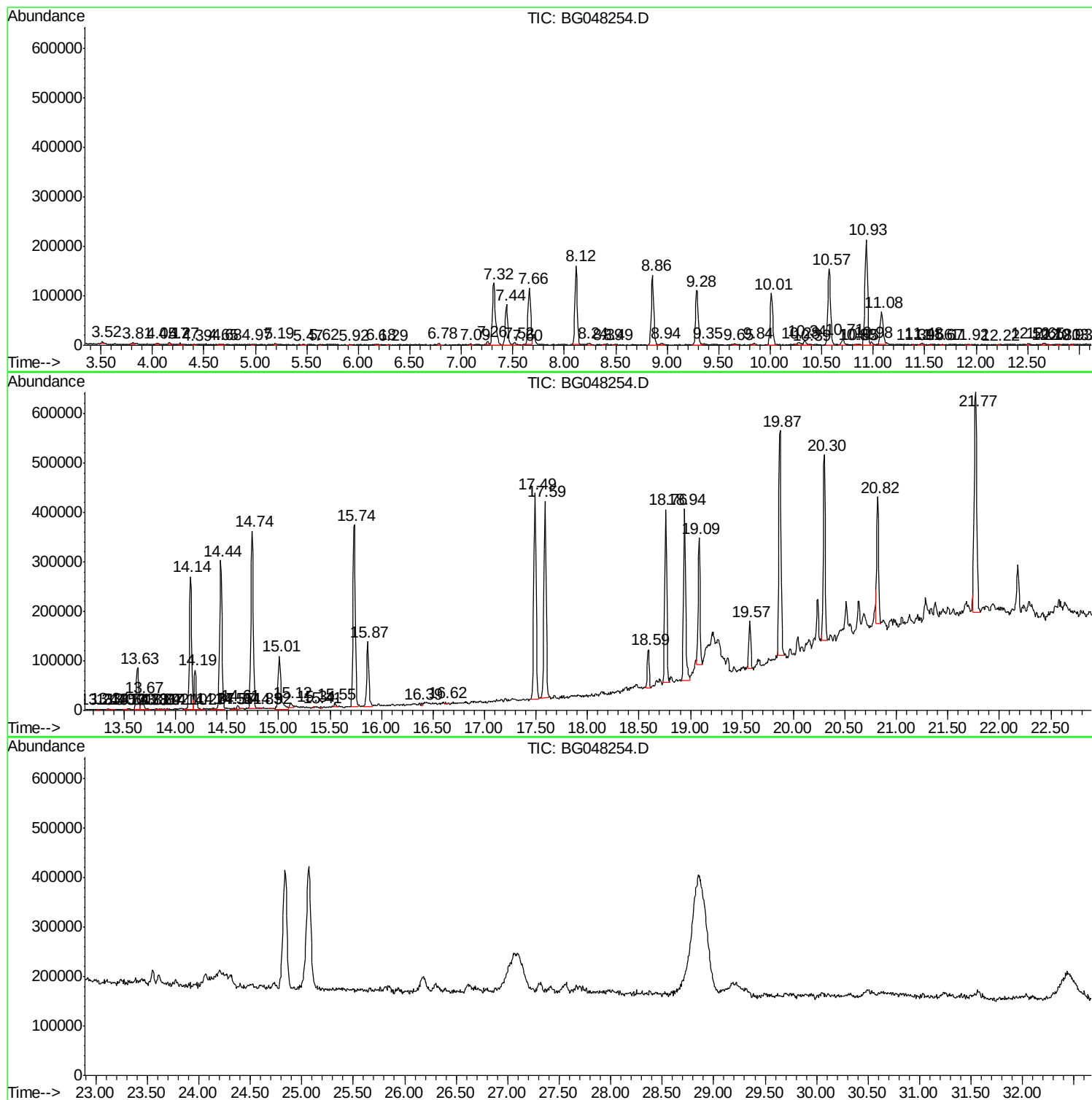
Sum of corrected areas: 10325577

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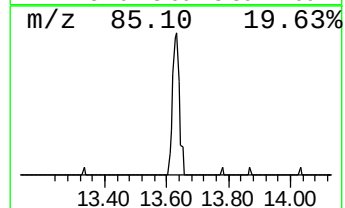
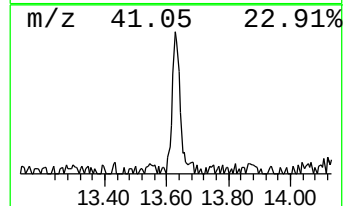
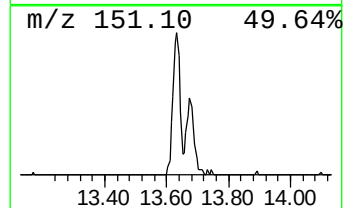
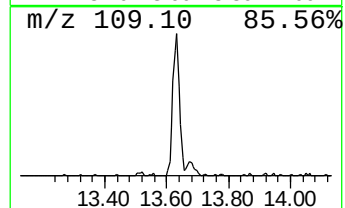
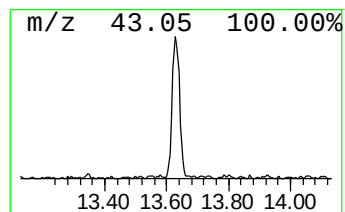
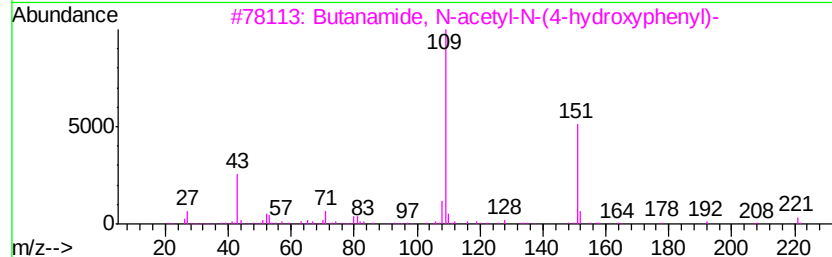
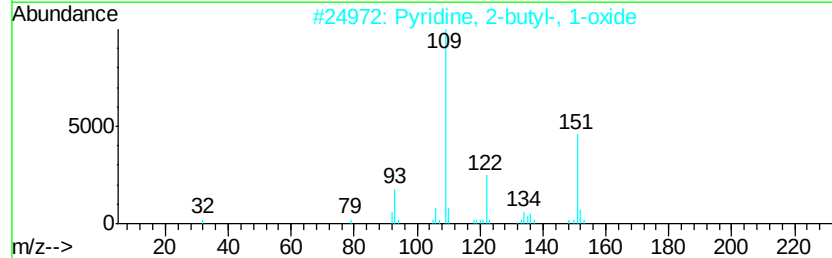
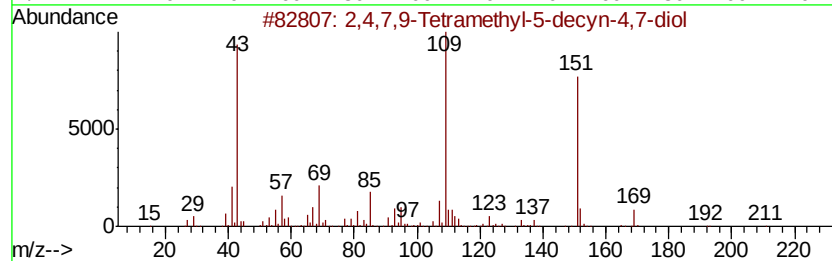
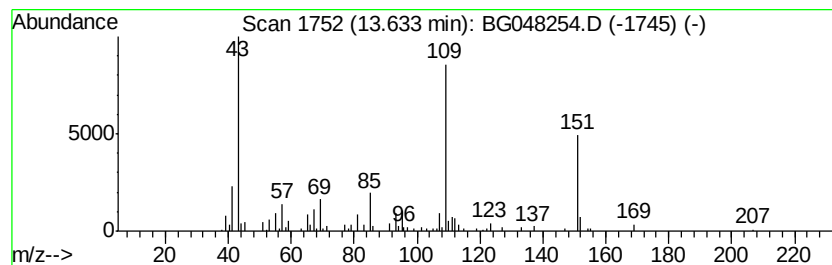
Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG021321MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 2,4,7,9-Tetramethyl-5-decyn... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.63	4.95 ng/ul	137741	Acenaphthene-d10	14.74	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4,7,9-Tetramethyl-5-decyn-4,7-...	226	C14H26O2	000126-86-3	91
2	Pyridine, 2-butyl-, 1-oxide	151	C9H13NO	031396-32-4	50
3	Butanamide, N-acetyl-N-(4-hydrox...	221	C12H15NO3	1000107-08-7	50
4	Metacetamol	151	C8H9NO2	000621-42-1	50
5	2-Propenal, 3-(2,2,6-trimethyl-7...	194	C12H18O2	055759-91-6	50



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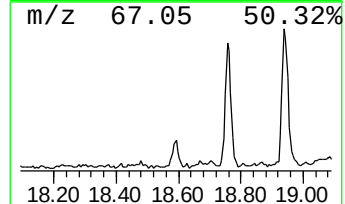
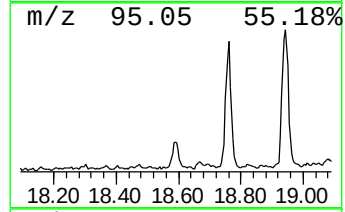
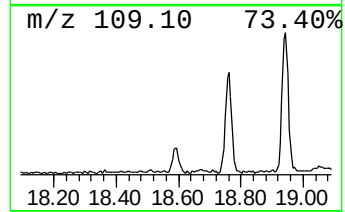
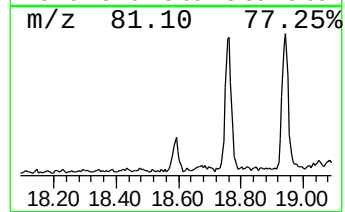
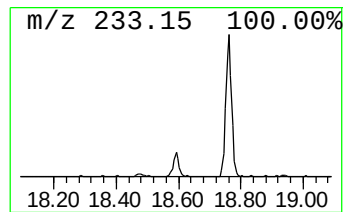
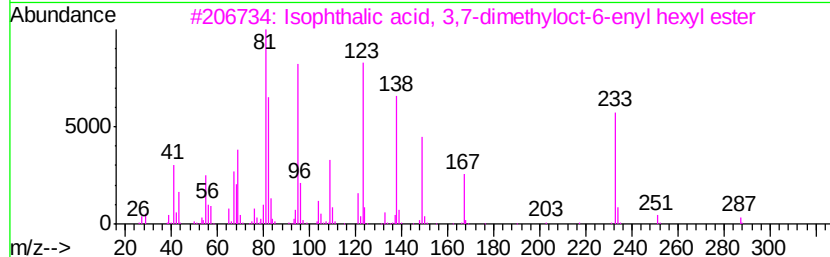
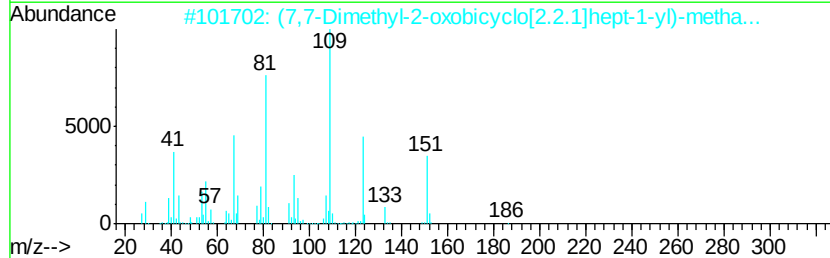
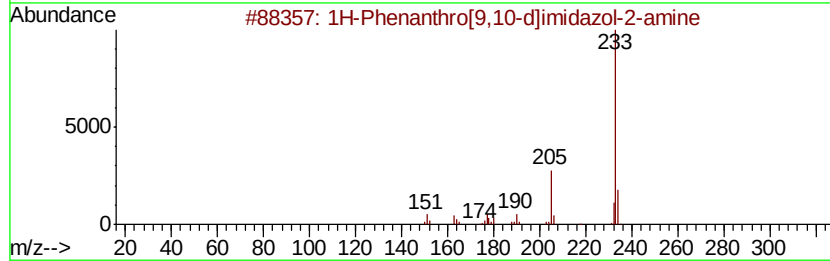
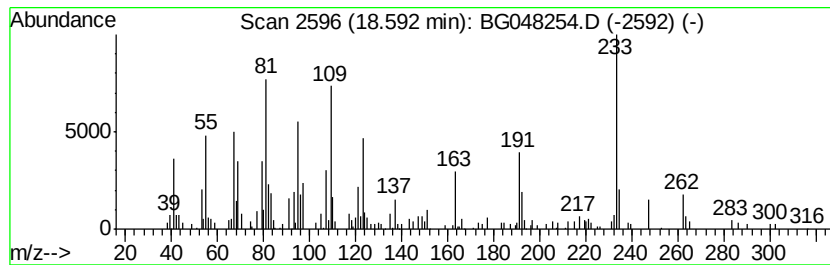
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Quant Title : SVOA CALIBRATION

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

Peak Number 2 unknown-01 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.		
18.59	3.12 ng/ul	102653	Phenanthrene-d10	17.49		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		1H-Phenanthro[9,10-d]imidazol-2-...	233	C15H11N3	037052-13-4	43
2		(7,7-Dimethyl-2-oxobicyclo[2.2.1]...	250	C10H15ClO3S	1000194-76-1	35
3		Isophthalic acid, 3,7-dimethyloc...	388	C24H36O4	1000356-74-0	35
4		18-Norabietane	262	C19H34	1000293-16-6	35
5		2-Cyclopenten-1-one, 3,4,4-trime...	124	C8H12O	030434-65-2	22



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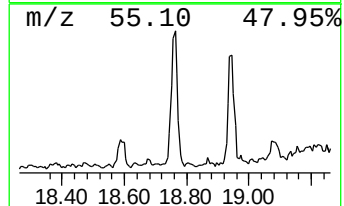
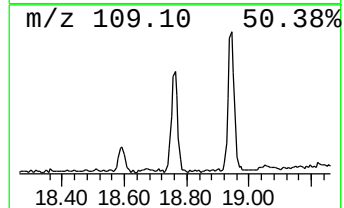
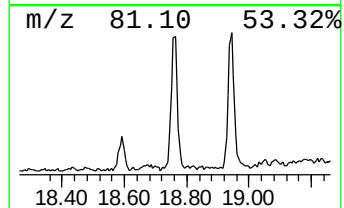
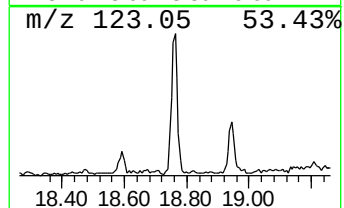
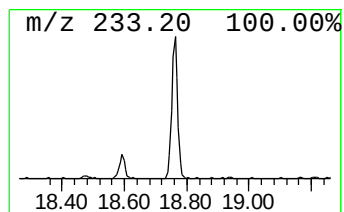
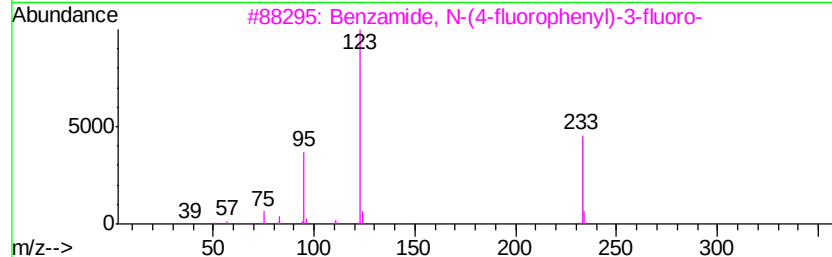
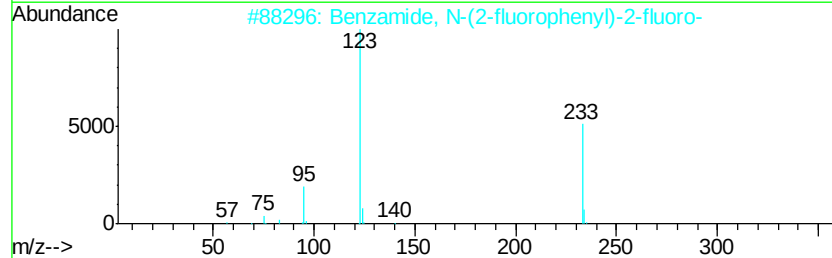
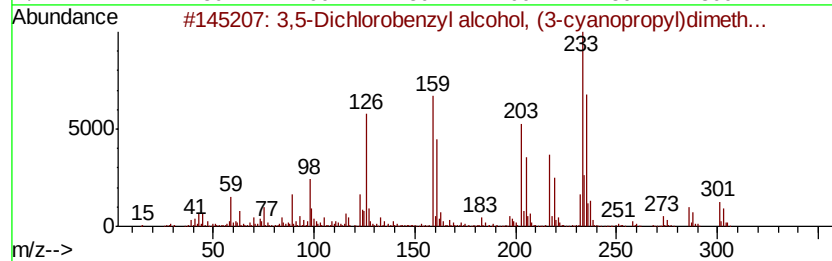
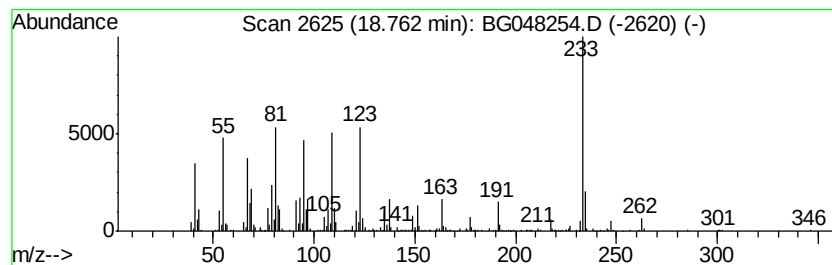
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TIC Integration Parameters: LSCINT.P

Peak Number 3 3,5-Dichlorobenzyl alcohol,... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.76	14.83 ng/ul	488170	Phenanthrene-d10	17.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3,5-Dichlorobenzyl alcohol, (3-c...	301	C13H17Cl2NOSi	1000376-10-3	86
2		Benzamide, N-(2-fluorophenyl)-2-...	233	C13H9F2NO	1000308-09-1	62
3		Benzamide, N-(4-fluorophenyl)-3-...	233	C13H9F2NO	1000307-16-3	58
4		18-Norabietane	262	C19H34	1000293-16-6	38
5		Benzamide, N-(4-fluorophenyl)-2-...	233	C13H9F2NO	1000307-07-9	38



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021921\
Data File : BG048254.D
Acq On : 21 Feb 2021 3:33
Operator : CG/JU
Sample : M1415-13
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBH27

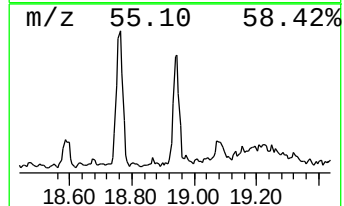
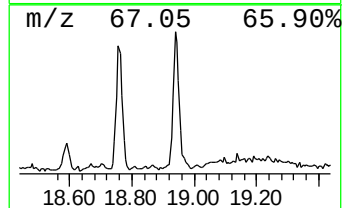
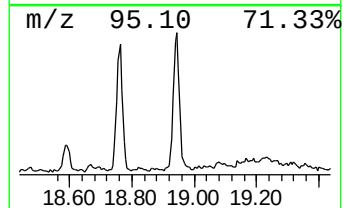
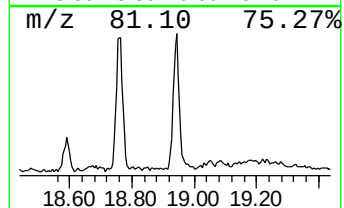
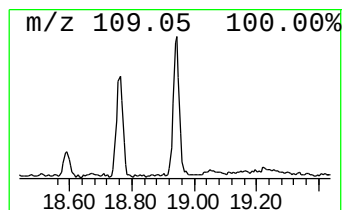
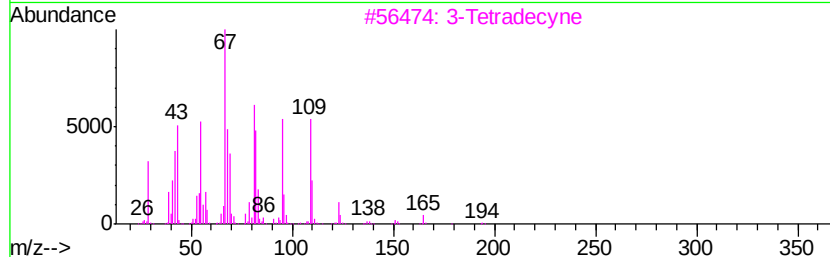
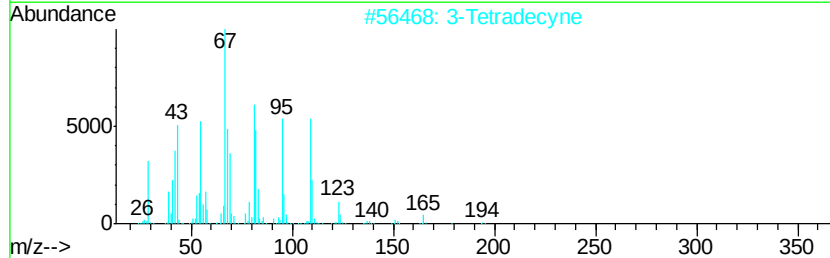
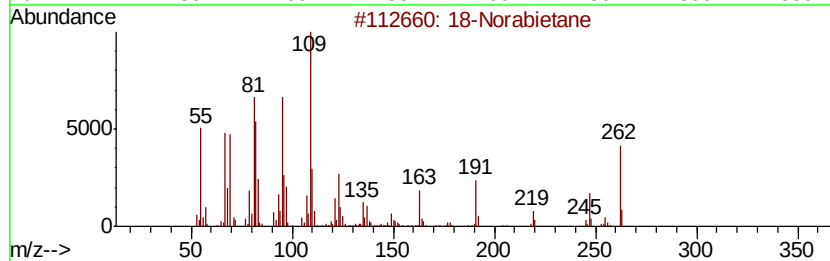
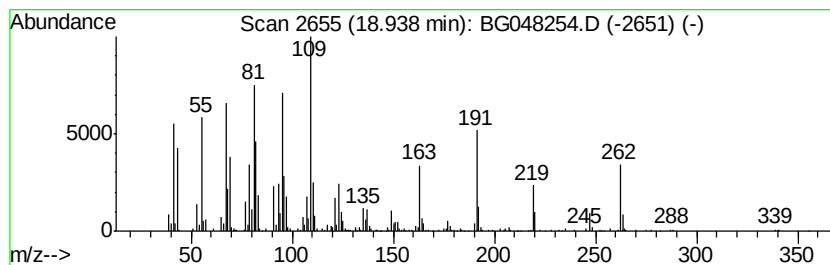
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 18-Norabietane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.94	15.15 ng/ul	498885	Phenanthrene-d10	17.49

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	18-Norabietane		262	C19H34	1000293-16-6	91
2	3-Tetradecyne		194	C14H26	060212-32-0	38
3	3-Tetradecyne		194	C14H26	060212-32-0	38
4	7-Octadecyne, 2-methyl-		264	C19H36	035354-38-2	35
5	7-Heptadecyne, 17-chloro-		270	C17H31Cl	056554-75-7	27



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021921\
Data File : BG048254.D
Acq On : 21 Feb 2021 3:33
Operator : CG/JU
Sample : M1415-13
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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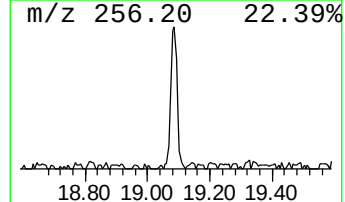
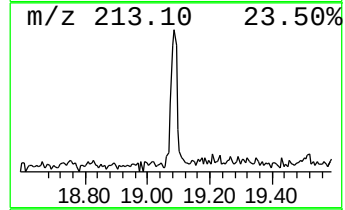
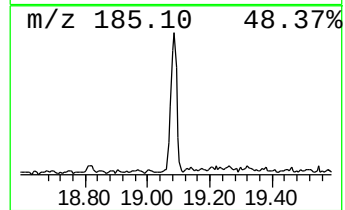
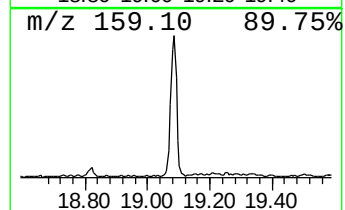
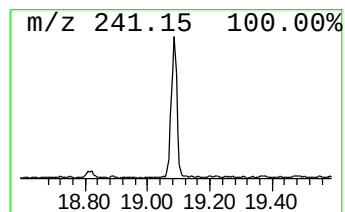
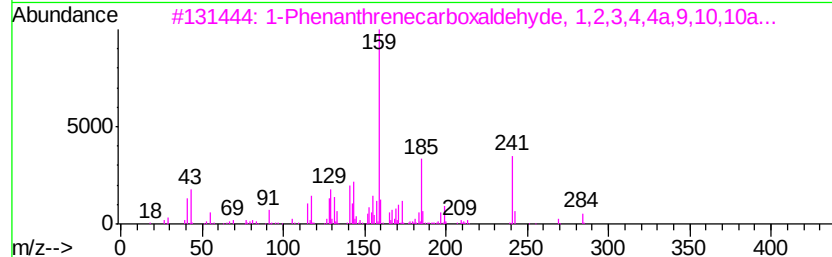
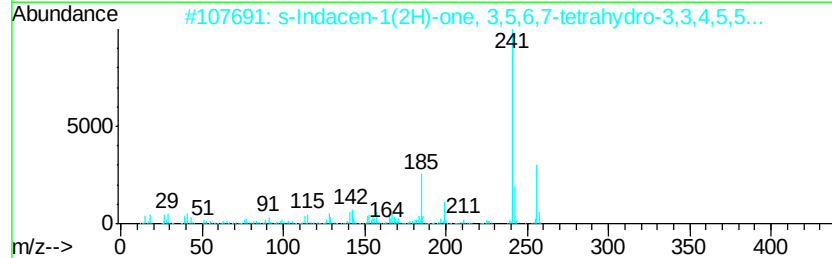
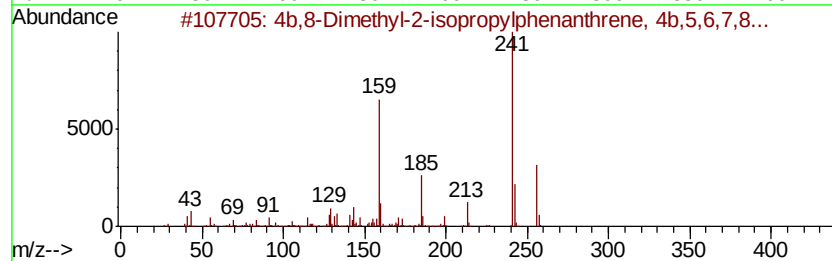
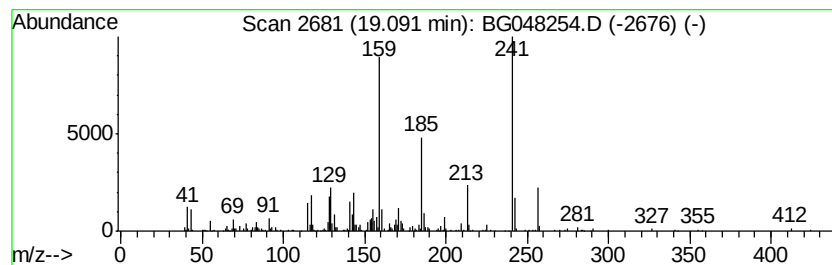
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 4b,8-Dimethyl-2-isopropylph... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.09	9.55 ng/ul	314489	Phenanthrene-d10	17.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4b,8-Dimethyl-2-isopropylphenant...	256	C19H28	1000197-14-1	97
2		s-Indacen-1(2H)-one, 3,5,6,7-tet...	256	C18H24O	038754-94-8	42
3		1-Phenanthrenecarboxaldehyde, 1,...	284	C20H28O	024035-50-5	38
4		3,4-Dihydro-3,3-dimethyl-1,9(2H,...	241	C15H15NO2	1000212-95-2	35
5		As-Indacene, 1,2,3,6,7,8-hexahyd...	256	C19H28	017465-47-3	30



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021921\
Data File : BG048254.D
Acq On : 21 Feb 2021 3:33
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Sample : M1415-13
Misc :
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Instrument :
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ClientSampleId :
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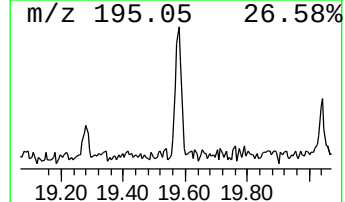
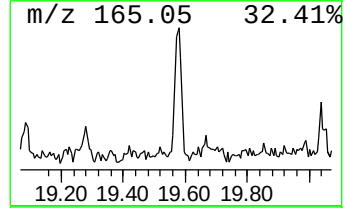
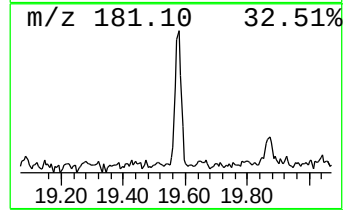
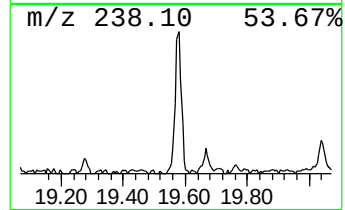
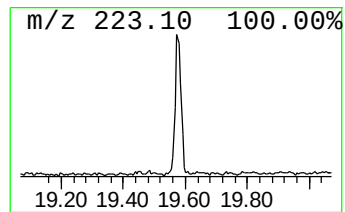
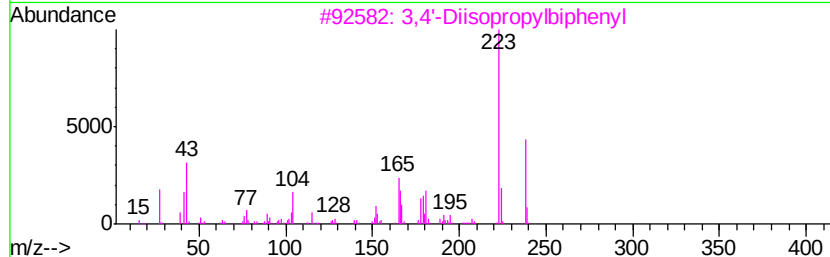
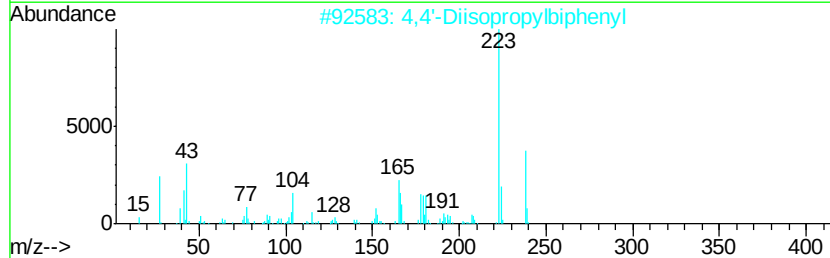
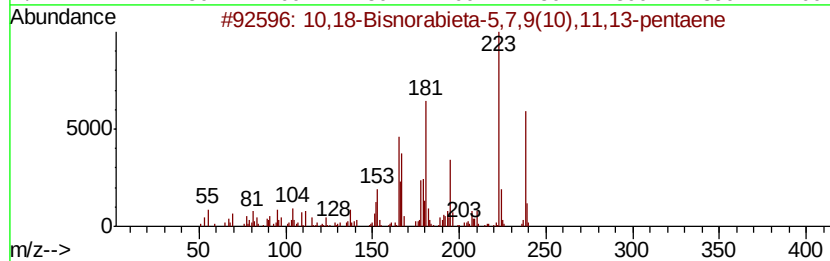
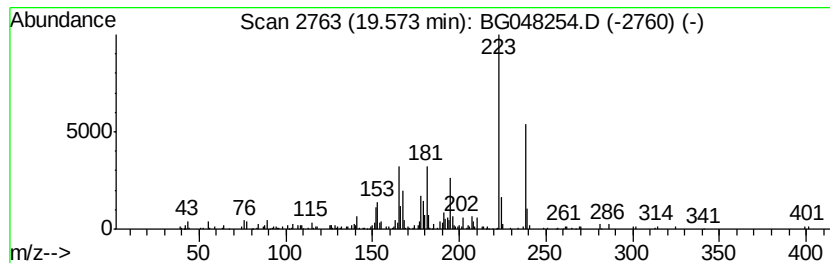
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 10,18-Bisnorabieta-5,7,9(10... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.57	3.86 ng/ul	127092	Phenanthrene-d10	17.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	10,18-Bisnorabieta-5,7,9(10),11,...	238	C18H22	006566-19-4	99
2		4,4'-Diisopropylbiphenyl	238	C18H22	018970-30-4	89
3		3,4'-Diisopropylbiphenyl	238	C18H22	061434-46-6	81
4		2,4(1H,3H)-Pteridinedione, 5,6,7...	238	C11H18N4O2	037921-31-6	64
5		3-(N,N-Diethylamino)carbazole	238	C16H18N2	054994-31-9	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021921\
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Operator : CG/JU
Sample : M1415-13
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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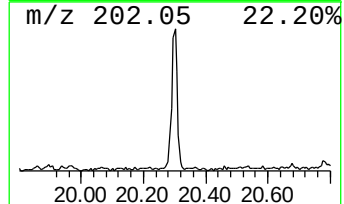
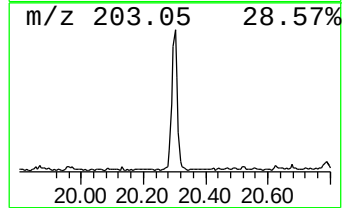
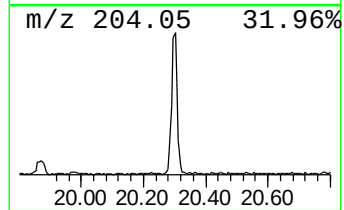
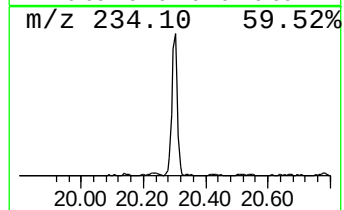
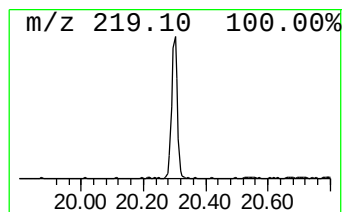
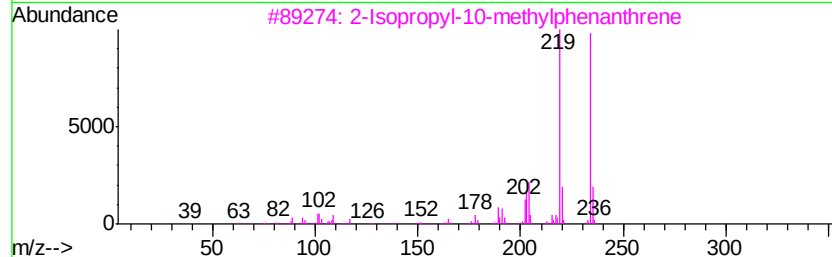
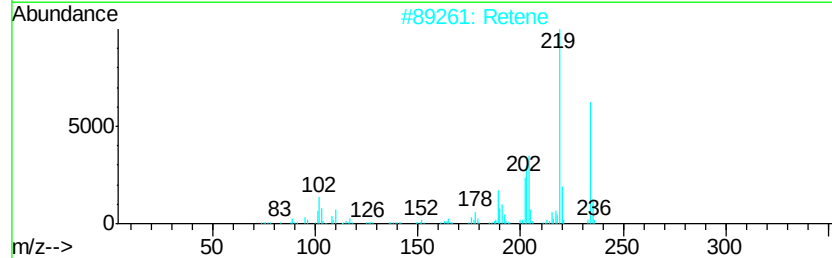
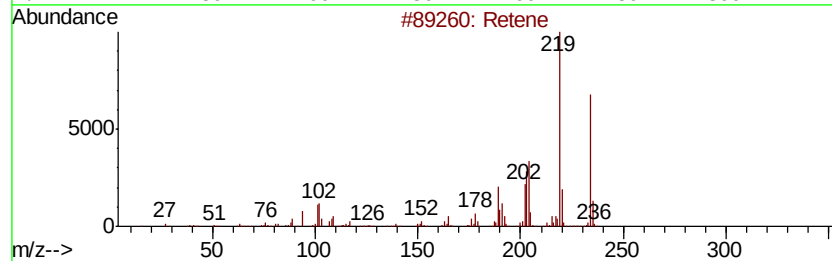
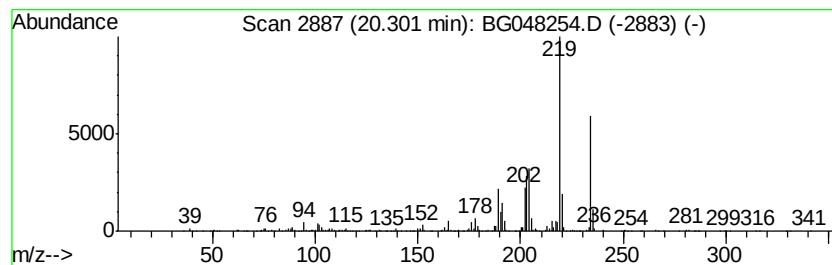
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Retene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.30	12.51 ng/ul	475189	Chrysene-d12	21.77

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	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021921\
Data File : BG048254.D
Acq On : 21 Feb 2021 3:33
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Sample : M1415-13
Misc :
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Instrument :
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ClientSampleId :
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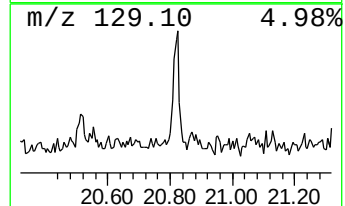
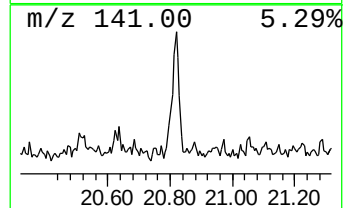
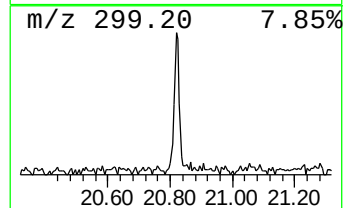
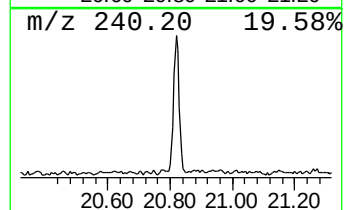
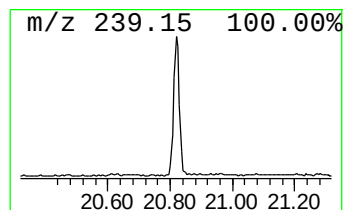
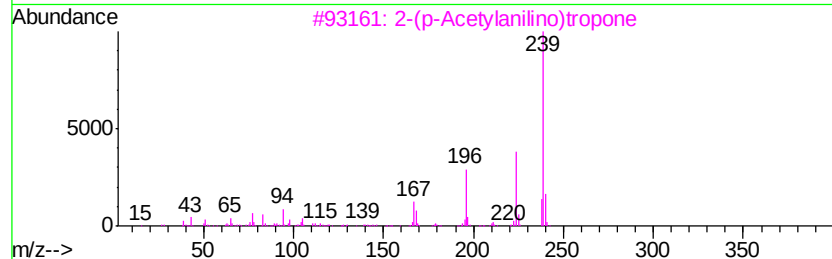
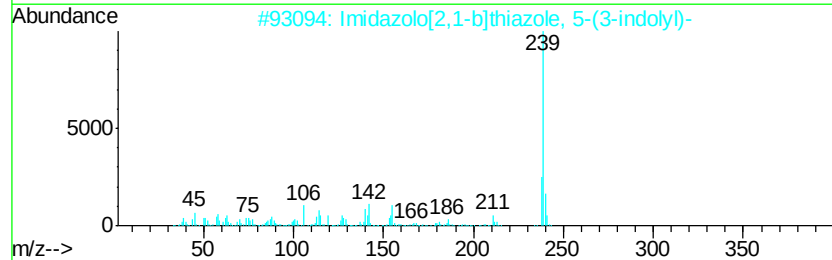
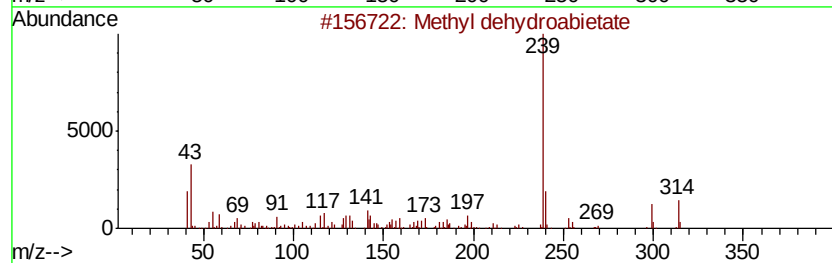
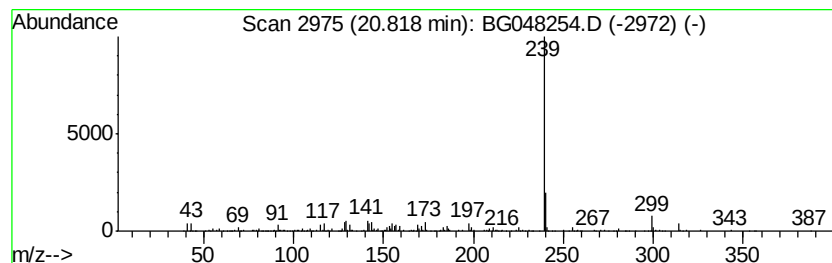
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Methyl dehydroabietate Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.82	8.56 ng/ul	325340	Chrysene-d12	21.77

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Methyl dehydroabietate	314	C21H30O2	001235-74-1	90
2		Imidazolo[2,1-b]thiazole, 5-(3-i...	239	C13H9N3S	327097-37-0	59
3		2-(p-Acetylanilino)tropone	239	C15H13NO2	1000241-52-9	50
4		trans-1,2-Bis(methyldichlorosily...	252	C4H8Cl4Si2	065899-10-7	47
5		9,10-Anthracenedione, 2-amino-3-...	239	C14H9NO3	000117-77-1	45



Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG021921\
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ALS Vial : 23 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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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
2,4,7,9-Tetrameth...	13.63	5.0	ng/ul	137741	3	14.74	556854	20.0
unknown-01	18.59	3.1	ng/ul	102653	4	17.49	658521	20.0
3,5-Dichlorobenzy...	18.76	14.8	ng/ul	488170	4	17.49	658521	20.0
18-Norabietane	18.94	15.2	ng/ul	498885	4	17.49	658521	20.0
4b,8-Dimethyl-2-i...	19.09	9.6	ng/ul	314489	4	17.49	658521	20.0
10,18-Bisnorabiet...	19.57	3.9	ng/ul	127092	4	17.49	658521	20.0
Retene	20.30	12.5	ng/ul	475189	5	21.77	759706	20.0
Methyl dehydroabi...	20.82	8.6	ng/ul	325340	5	21.77	759706	20.0