

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111324\
 Data File : BG063413.D
 Acq On : 15 Nov 2024 5:25
 Operator : RC/JU
 Sample : P4761-03
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 A0AL4

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: BG063413.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.300	31	36	39	rBV5	26645	34715	0.88%	0.072%
2	3.341	39	43	48	rVB4	33830	48885	1.24%	0.101%
3	3.576	79	83	86	rVB4	11425	14016	0.35%	0.029%
4	3.711	101	106	114	rBV	264057	450546	11.40%	0.934%
5	4.228	190	194	199	rVB4	6815	12269	0.31%	0.025%
6	4.486	233	238	243	rVV6	14095	25274	0.64%	0.052%
7	4.945	312	316	319	rBV5	10986	15882	0.40%	0.033%
8	4.992	322	324	328	rVV4	8891	13166	0.33%	0.027%
9	5.397	390	393	395	rVV3	8572	10693	0.27%	0.022%
10	5.491	404	409	416	rVV8	11124	21012	0.53%	0.044%
11	5.538	416	417	424	rVB4	7672	12643	0.32%	0.026%
12	5.932	478	484	489	rVB8	11730	22968	0.58%	0.048%
13	6.108	508	514	515	rBV6	10131	14921	0.38%	0.031%
14	6.255	535	539	542	rBV5	5213	9572	0.24%	0.020%
15	6.337	547	553	556	rBV8	10184	20972	0.53%	0.043%
16	6.508	579	582	588	rVB6	14186	16691	0.42%	0.035%
17	7.007	657	667	679	rBV	208759	402338	10.18%	0.834%
18	7.166	688	694	701	rBV	378453	640217	16.21%	1.327%
19	7.365	719	728	737	rBV	486721	856624	21.68%	1.776%
20	7.430	737	739	743	rVV5	8685	10614	0.27%	0.022%
21	7.830	800	807	813	rBV2	468386	779848	19.74%	1.616%
22	7.900	813	819	828	rVB	605834	1000299	25.32%	2.073%
23	8.159	858	863	864	rVV5	7717	11502	0.29%	0.024%
24	8.176	864	866	867	rVV2	11021	9849	0.25%	0.020%
25	8.211	867	872	878	rVB3	26869	46571	1.18%	0.097%
26	8.482	913	918	920	rVB5	9398	14290	0.36%	0.030%
27	8.541	922	928	938	rVV3	557899	992819	25.13%	2.058%
28	8.611	938	940	945	rVV6	12058	17676	0.45%	0.037%
29	8.658	945	948	952	rVB5	8591	11617	0.29%	0.024%
30	8.811	966	974	976	rBV5	23096	47366	1.20%	0.098%
31	8.934	985	995	998	rBV6	57938	119796	3.03%	0.248%
32	8.987	998	1004	1012	rVV2	497507	890687	22.55%	1.846%
33	9.175	1031	1036	1042	rVV9	18368	45420	1.15%	0.094%
34	9.287	1050	1055	1062	rVV	168160	270510	6.85%	0.561%
35	9.334	1062	1063	1069	rVB6	6923	10574	0.27%	0.022%
36	9.457	1077	1084	1089	rBV3	19764	28211	0.71%	0.058%

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

37	9.710	1119	1127	1134	rBV2	471820	852011	21.57%	1.766%
38	9.810	1138	1144	1146	rBV7	16966	31309	0.79%	0.065%
39	9.898	1154	1159	1168	rVB	314632	522251	13.22%	1.082%
40	10.039	1174	1183	1190	rBV4	92089	215970	5.47%	0.448%
41	10.103	1193	1194	1199	rVB5	11376	9480	0.24%	0.020%
42	10.250	1212	1219	1224	rBV	714687	1291566	32.69%	2.677%
43	10.291	1224	1226	1233	rVB2	133409	199883	5.06%	0.414%
44	10.620	1275	1282	1290	rBV2	657489	1128000	28.55%	2.338%
45	10.761	1298	1306	1310	rBV	407865	780928	19.77%	1.619%
46	10.797	1310	1312	1320	rVB2	123740	187321	4.74%	0.388%
47	11.085	1358	1361	1363	rVV4	9130	11492	0.29%	0.024%
48	11.214	1380	1383	1386	rBV5	17645	21384	0.54%	0.044%
49	11.378	1406	1411	1418	rBV6	37276	97142	2.46%	0.201%
50	11.502	1426	1432	1436	rBV6	37272	89589	2.27%	0.186%
51	11.584	1441	1446	1452	rVB	92571	161338	4.08%	0.334%
52	11.825	1483	1487	1489	rBV5	27354	44194	1.12%	0.092%
53	12.119	1529	1537	1542	rBV5	23239	56017	1.42%	0.116%
54	12.195	1544	1550	1555	rBV9	42744	90141	2.28%	0.187%
55	12.348	1572	1576	1577	rBV4	11397	15248	0.39%	0.032%
56	12.383	1580	1582	1584	rBV3	14368	13871	0.35%	0.029%
57	12.559	1608	1612	1615	rBV4	35841	55127	1.40%	0.114%
58	12.694	1633	1635	1636	rBV2	17008	15831	0.40%	0.033%
59	12.806	1651	1654	1657	rBV5	16434	25317	0.64%	0.052%
60	12.871	1657	1665	1674	rVB	519458	835823	21.16%	1.732%
61	12.941	1675	1677	1680	rBV4	10996	12461	0.32%	0.026%
62	12.994	1680	1686	1696	rVB4	30965	91711	2.32%	0.190%
63	13.188	1715	1719	1720	rBV4	18476	21796	0.55%	0.045%
64	13.282	1732	1735	1736	rBV3	16678	17071	0.43%	0.035%
65	13.411	1754	1757	1758	rBV3	16274	17358	0.44%	0.036%
66	13.505	1769	1773	1782	rBV3	20806	61164	1.55%	0.127%
67	13.682	1799	1803	1805	rBV5	15623	18994	0.48%	0.039%
68	13.875	1830	1836	1846	rVB	1346103	1959338	49.60%	4.061%
69	13.999	1851	1857	1865	rVB	301678	450538	11.40%	0.934%
70	14.075	1865	1870	1871	rBV5	24362	37065	0.94%	0.077%
71	14.116	1872	1877	1879	rVV4	86736	174384	4.41%	0.361%
72	14.163	1879	1885	1897	rVB	1426877	2235751	56.59%	4.634%
73	14.304	1904	1909	1918	rBV2	110938	261935	6.63%	0.543%
74	14.392	1921	1924	1930	rVV8	53011	91220	2.31%	0.189%
75	14.469	1932	1937	1946	rVV	1202163	1904823	48.22%	3.948%
76	14.533	1946	1948	1953	rVB3	44012	51401	1.30%	0.107%

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 Title : SVOA CALIBRATION

77	14.698	1970	1976	1981	rBV3	155739	299485	7.58%	0.621%
78	14.874	2003	2006	2010	rBV5	40924	65345	1.65%	0.135%
79	15.215	2062	2064	2066	rBV3	38770	41846	1.06%	0.087%
80	15.244	2067	2069	2075	rVB7	66120	106162	2.69%	0.220%
81	15.468	2101	2107	2113	rBV	1982019	2917062	73.84%	6.046%
82	15.521	2114	2116	2120	rVB5	68488	77421	1.96%	0.160%
83	15.591	2123	2128	2138	rBV	708887	1093558	27.68%	2.267%
84	15.691	2141	2145	2148	rBV2	95655	117758	2.98%	0.244%
85	15.832	2164	2169	2176	rVB	247241	375451	9.50%	0.778%
86	15.979	2189	2194	2199	rBV3	137565	244770	6.20%	0.507%
87	16.396	2263	2265	2269	rVB5	39924	41525	1.05%	0.086%
88	16.490	2278	2281	2286	rBV2	62010	74443	1.88%	0.154%
89	17.224	2400	2406	2413	rVB	1484237	2272824	57.53%	4.711%
90	17.324	2417	2423	2430	rVB	2105685	3205881	81.15%	6.645%
91	17.389	2432	2434	2437	rBV3	36788	42574	1.08%	0.088%
92	17.718	2486	2490	2499	rBV	238285	396480	10.04%	0.822%
93	18.129	2556	2560	2566	rBV5	143009	209533	5.30%	0.434%
94	18.341	2592	2596	2603	rBV	207561	361233	9.14%	0.749%
95	18.505	2620	2624	2629	rBV4	149655	258152	6.53%	0.535%
96	19.622	2808	2814	2818	rBV	2925730	3950477	100.00%	8.188%
97	19.669	2818	2822	2832	rVB	1259083	1788196	45.27%	3.706%
98	21.478	3124	3130	3138	rVB2	1836199	2708720	68.57%	5.614%
99	24.310	3602	3612	3624	rVB	1449454	3739030	94.65%	7.750%
100	24.516	3638	3647	3657	rVB	1063956	2748556	69.58%	5.697%

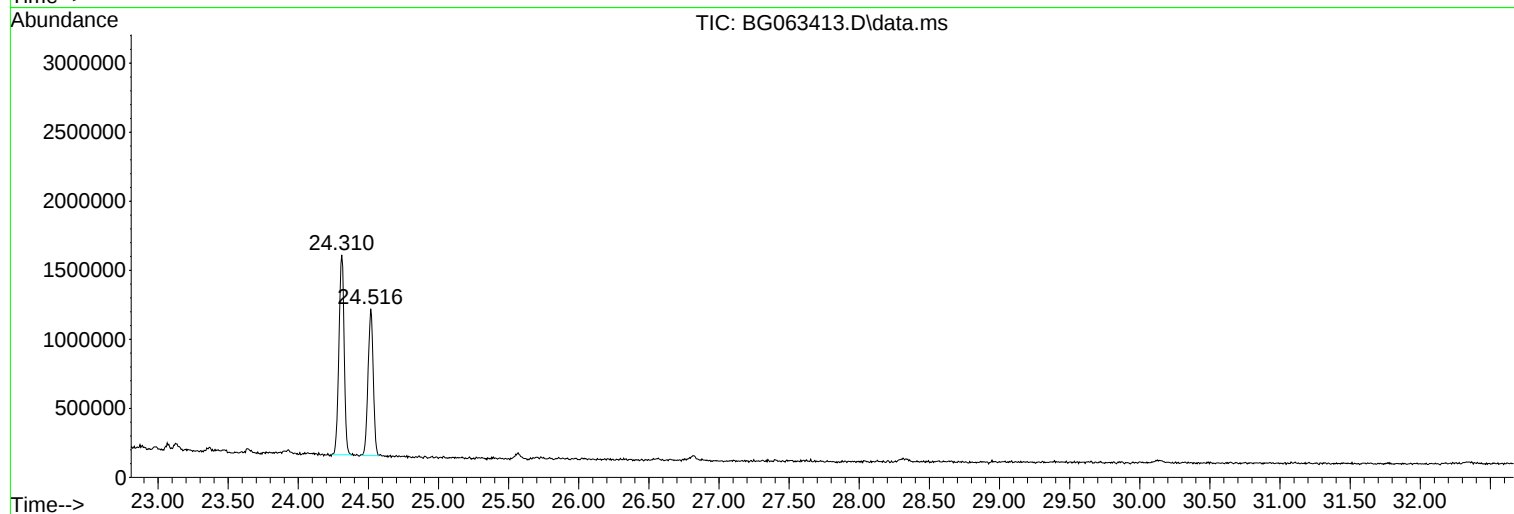
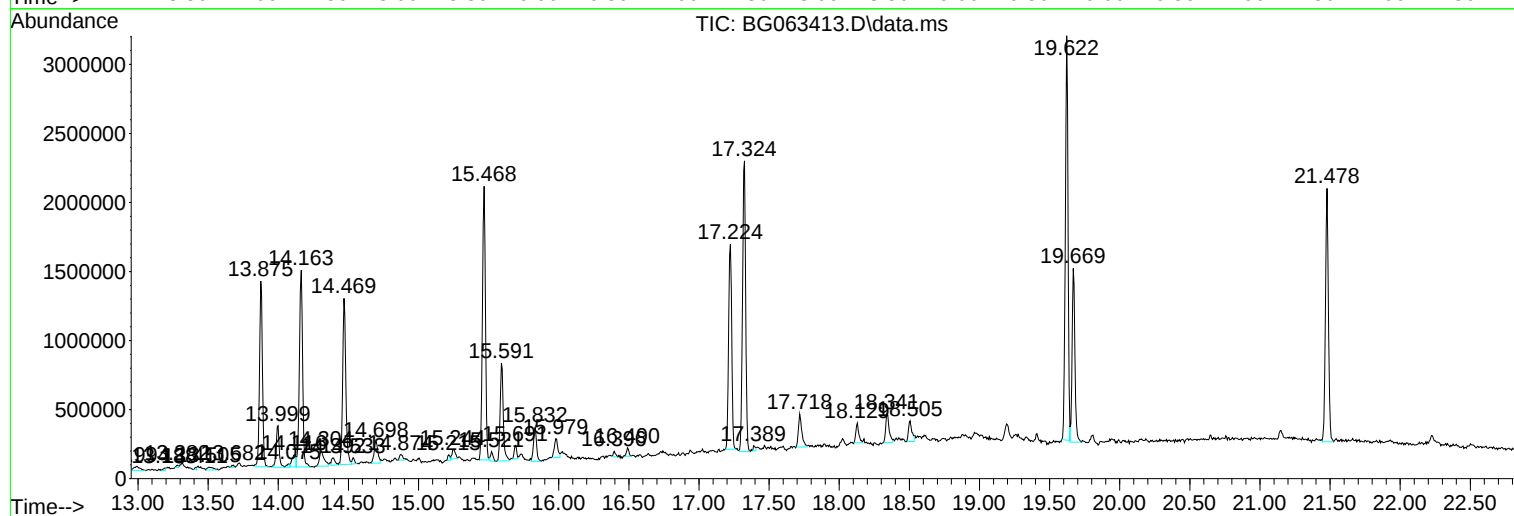
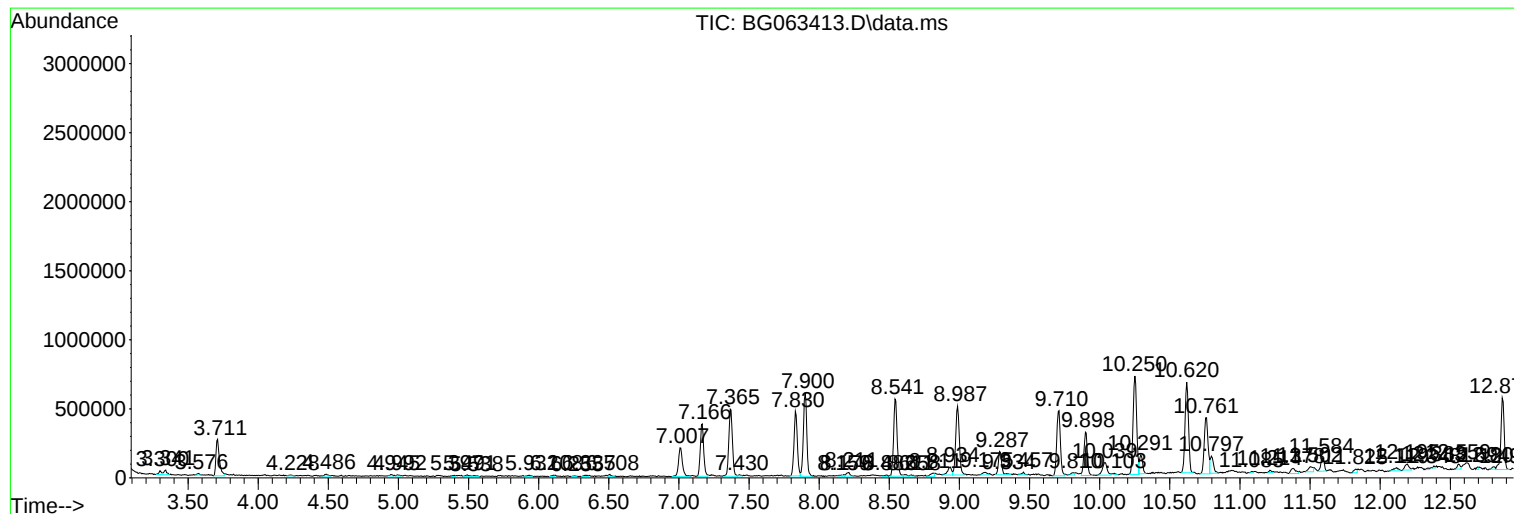
Sum of corrected areas: 48245778

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Misc :
ALS Vial : 29 Sample Multiplier: 1

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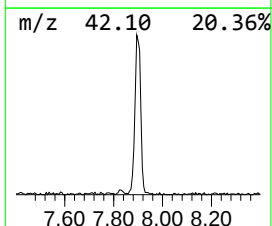
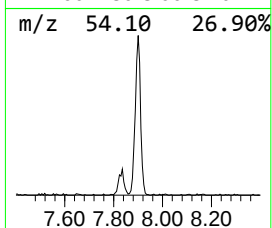
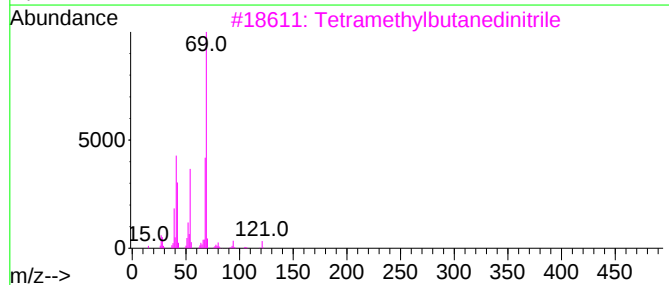
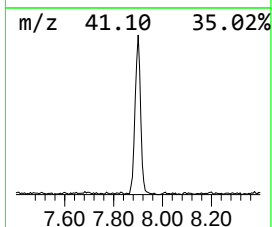
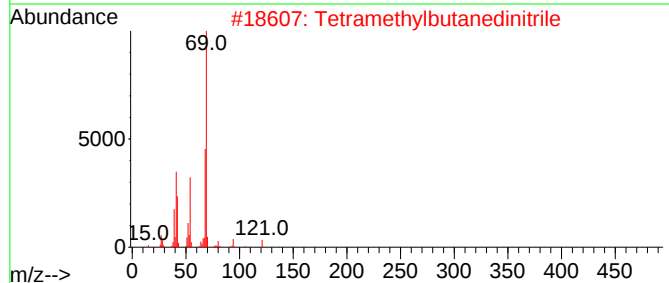
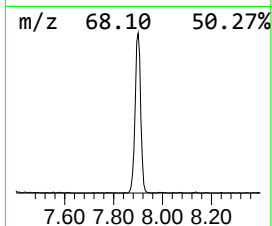
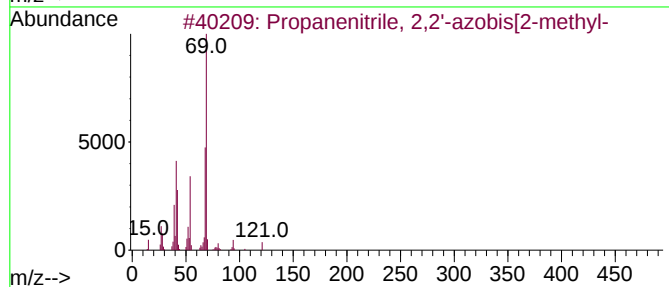
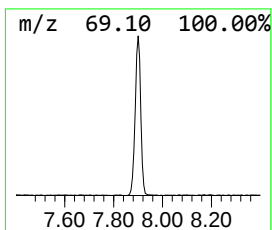
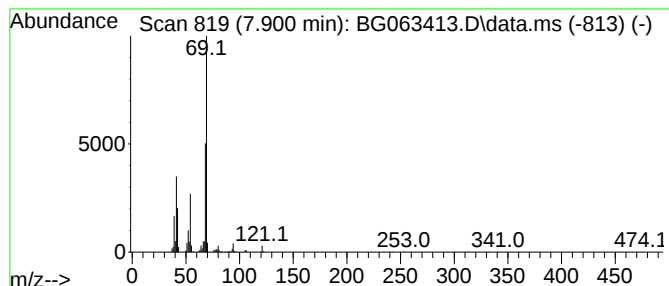
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110624.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Propanenitrile, 2,2'-azobis... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.900	25.65 ng/ul	1000300	1,4-Dichlorobenzene-d4	7.830

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propanenitrile, 2,2'-azobis[2-me...	164	C8H12N4	000078-67-1	90
2			Tetramethylbutanedinitrile	136	C8H12N2	003333-52-6	90
3			Tetramethylbutanedinitrile	136	C8H12N2	003333-52-6	86
4			Propanenitrile, 2,2'-azobis[2-me...	164	C8H12N4	000078-67-1	78
5			Propanenitrile, 2,2'-azobis[2-me...	164	C8H12N4	000078-67-1	72



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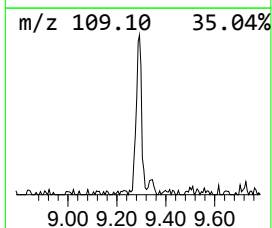
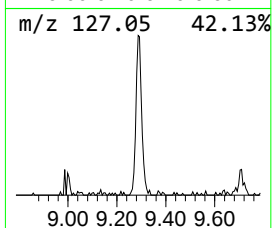
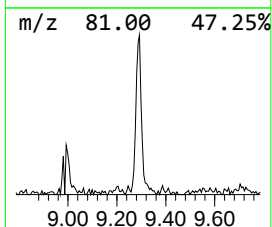
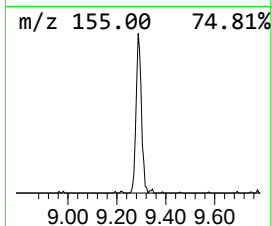
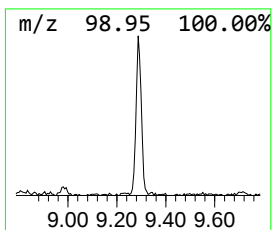
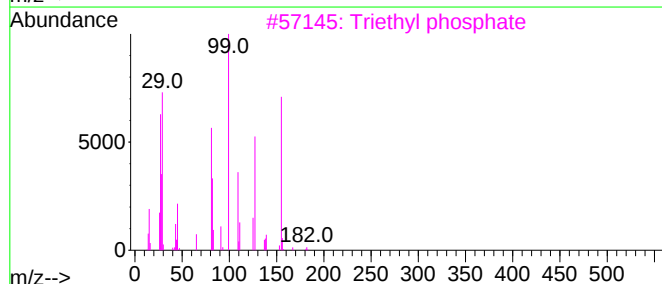
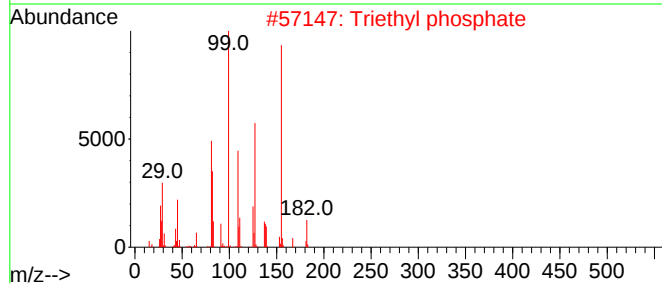
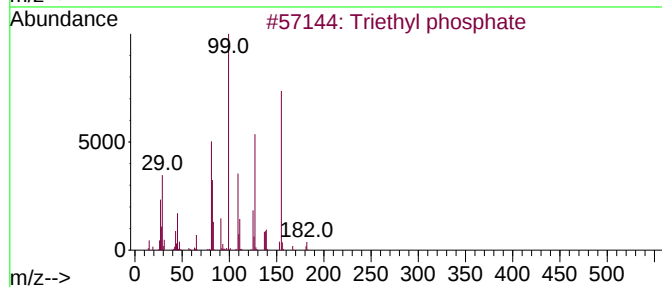
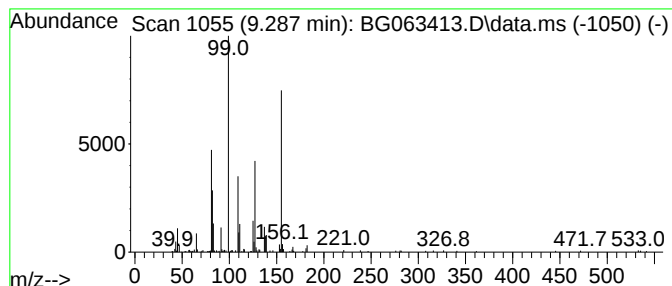
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110624.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Triethyl phosphate Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.287	4.80 ng/ul	270510	Naphthalene-d8	10.620

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Triethyl phosphate	182	C6H15O4P	000078-40-0	98
2			Triethyl phosphate	182	C6H15O4P	000078-40-0	97
3			Triethyl phosphate	182	C6H15O4P	000078-40-0	78
4			Butyl diethyl phosphate	210	C8H19O4P	1000298-58-0	64
5			Triethyl phosphate	182	C6H15O4P	000078-40-0	64



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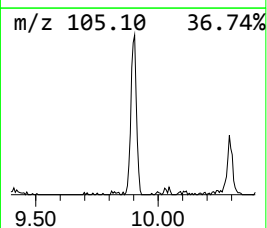
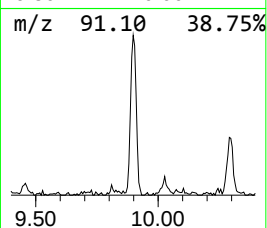
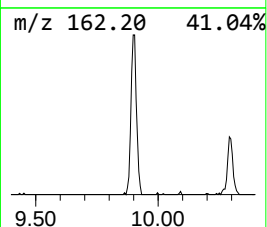
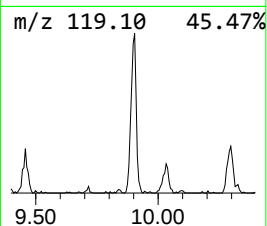
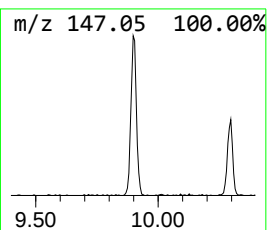
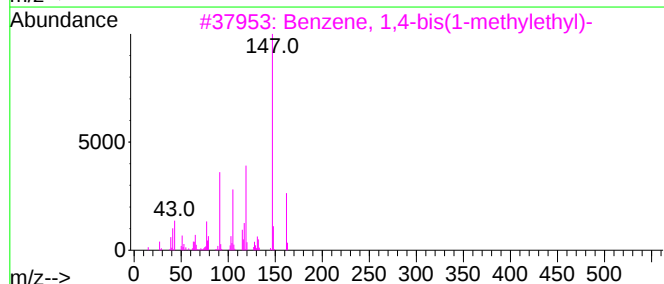
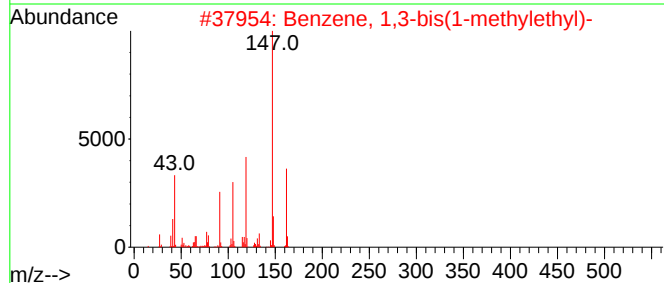
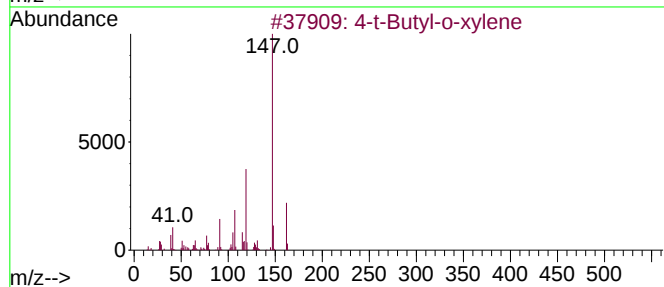
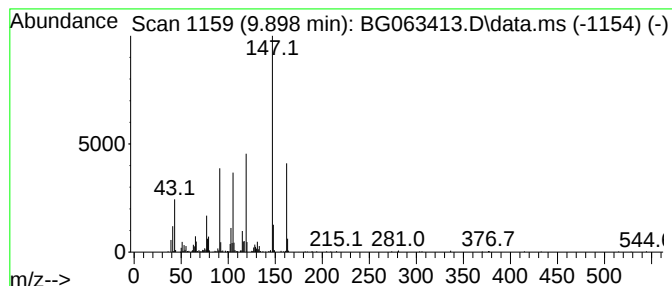
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 4-t-Butyl-o-xylene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.898	9.26 ng/ul	522251	Naphthalene-d8	10.620

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-t-Butyl-o-xylene	162	C12H18	007397-06-0	93
2			Benzene, 1,3-bis(1-methylethyl)-	162	C12H18	000099-62-7	93
3			Benzene, 1,4-bis(1-methylethyl)-	162	C12H18	000100-18-5	91
4			Benzene, 1,3-bis(1-methylethyl)-	162	C12H18	000099-62-7	91
5			Benzene, 1,2-bis(1-methylethyl)-	162	C12H18	000577-55-9	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111324\
Data File : BG063413.D
Acq On : 15 Nov 2024 5:25
Operator : RC/JU
Sample : P4761-03
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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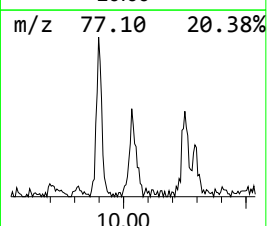
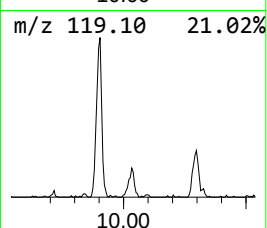
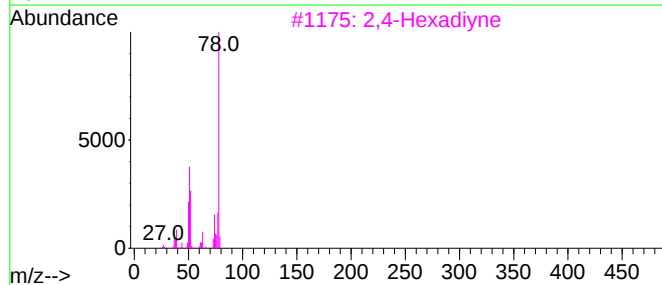
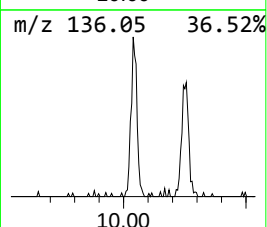
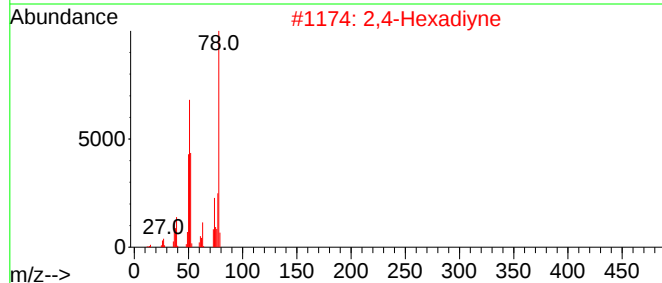
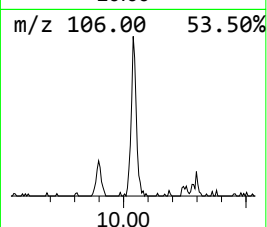
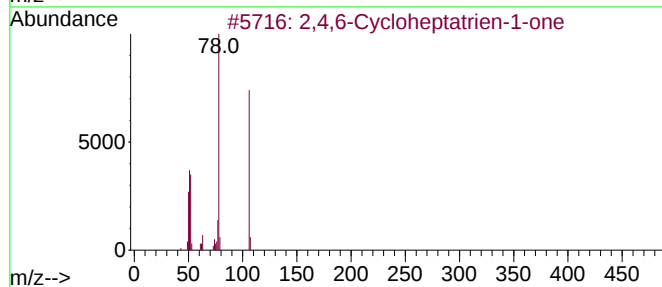
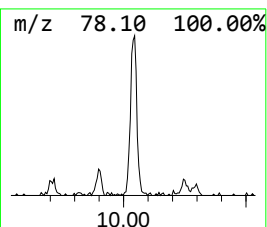
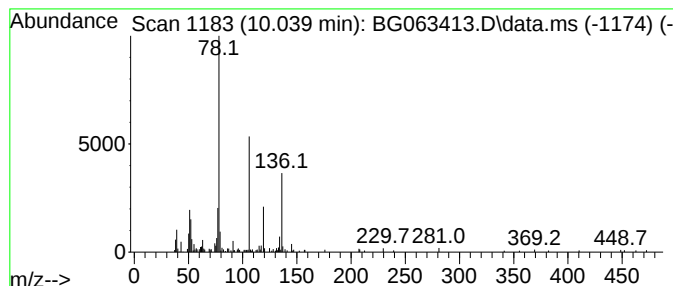
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 2,4,6-Cycloheptatrien-1-one Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.039	3.83 ng/ul	215970	Naphthalene-d8	10.620

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2,4,6-Cycloheptatrien-1-one	106	C7H6O	000539-80-0	87
2		2,4-Hexadiyne	78	C6H6	002809-69-0	81
3		2,4-Hexadiyne	78	C6H6	002809-69-0	74
4		4H-1,3-Benzodioxin	136	C8H8O2	000254-27-3	72
5		2,4-Hexadiyne	78	C6H6	002809-69-0	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111324\
Data File : BG063413.D
Acq On : 15 Nov 2024 5:25
Operator : RC/JU
Sample : P4761-03
Misc :
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Instrument :
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ClientSampleId :
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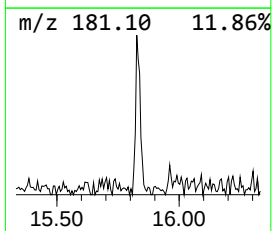
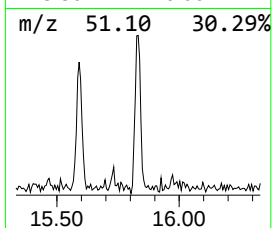
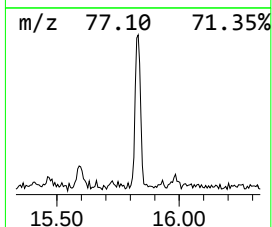
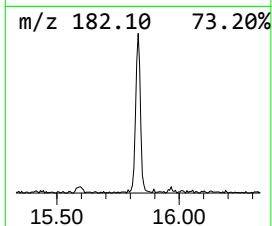
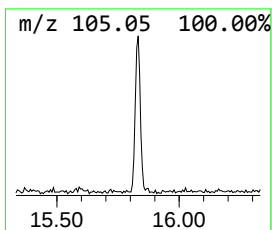
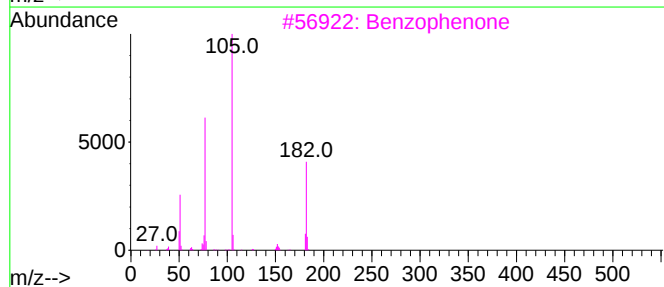
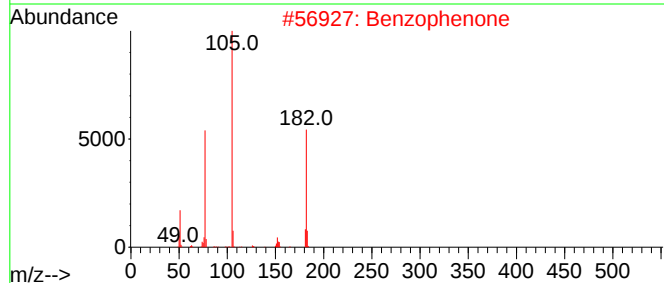
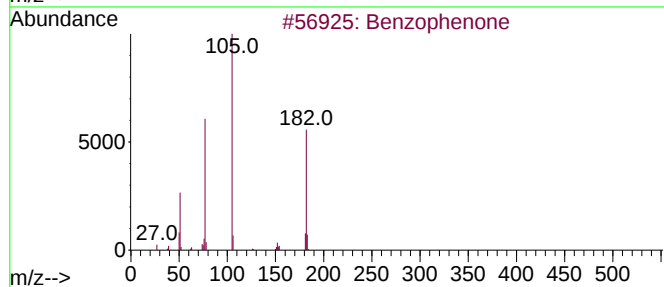
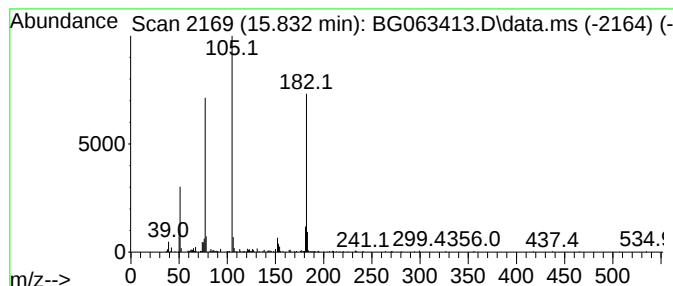
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Benzophenone Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.832	3.94 ng/ul	375451	Acenaphthene-d10	14.469

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	97
2			Benzophenone	182	C13H10O	000119-61-9	95
3			Benzophenone	182	C13H10O	000119-61-9	94
4			Benzophenone	182	C13H10O	000119-61-9	91
5			Benzophenone	182	C13H10O	000119-61-9	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111324\
Data File : BG063413.D
Acq On : 15 Nov 2024 5:25
Operator : RC/JU
Sample : P4761-03
Misc :
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Instrument :
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ClientSampleId :
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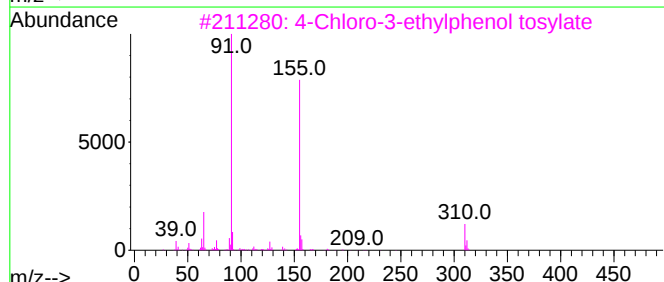
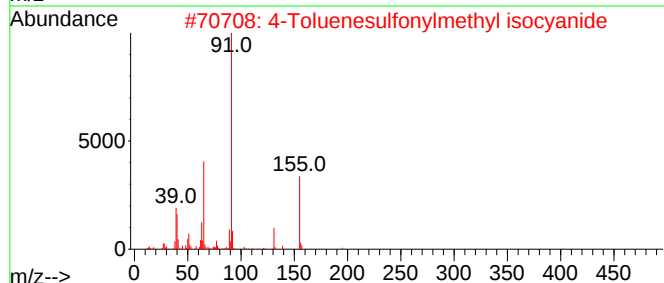
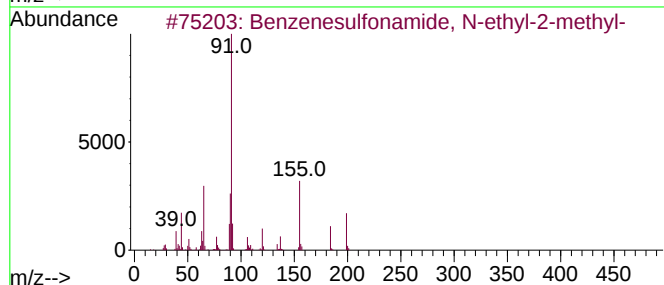
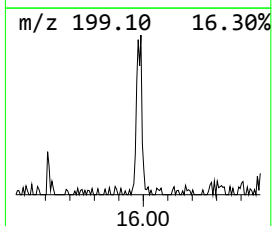
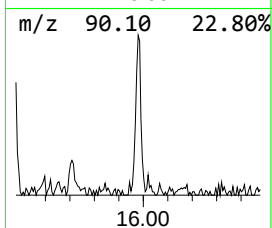
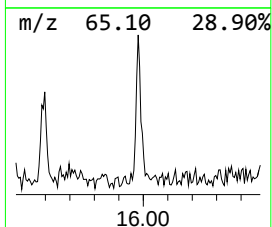
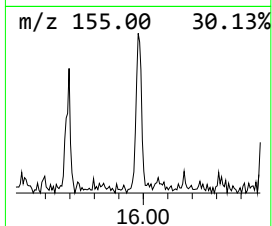
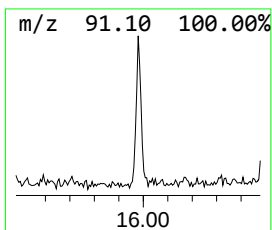
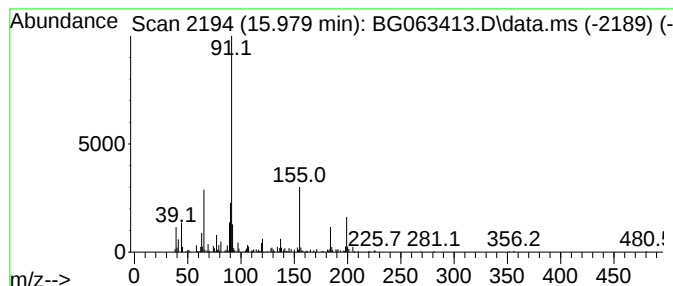
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 Benzenesulfonamide, N-ethyl... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.979	2.15 ng/ul	244770	Phenanthrene-d10	17.224

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzenesulfonamide, N-ethyl-2-me...	199	C9H13NO2S	001077-56-1	91
2			4-Toluenesulfonylmethyl isocyanide	195	C9H9NO2S	036635-61-7	50
3			4-Chloro-3-ethylphenol tosylate	310	C15H15ClO3S	1000467-53-5	47
4			Benzenesulfonic acid, 3-methyl-,...	200	C9H12O3S	126865-08-5	47
5			Benzenesulfonamide, N-ethyl-4-me...	199	C9H13NO2S	000080-39-7	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111324\
Data File : BG063413.D
Acq On : 15 Nov 2024 5:25
Operator : RC/JU
Sample : P4761-03
Misc :
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Instrument :
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ClientSampleId :
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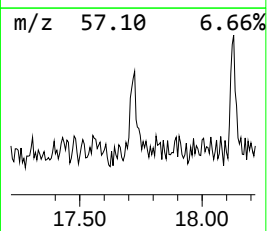
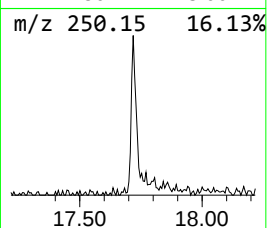
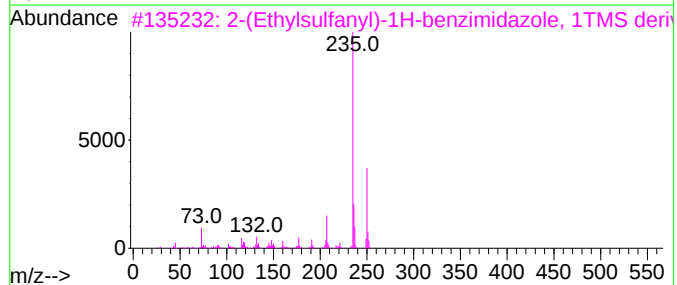
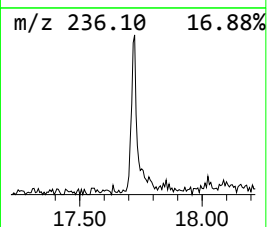
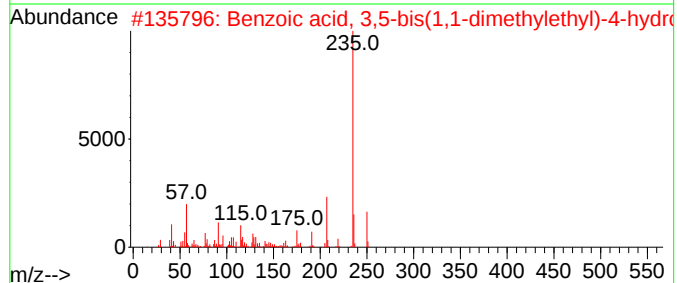
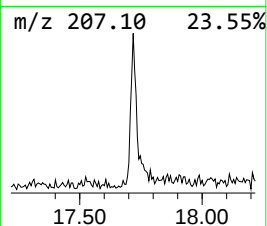
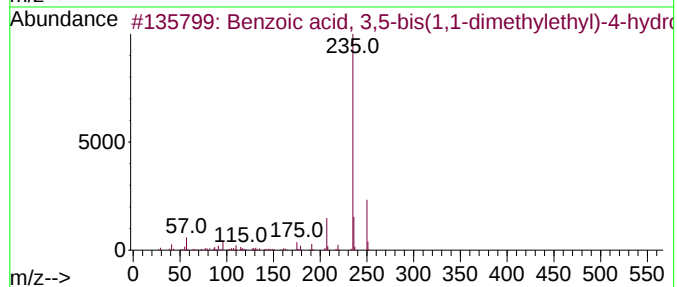
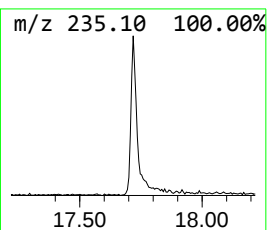
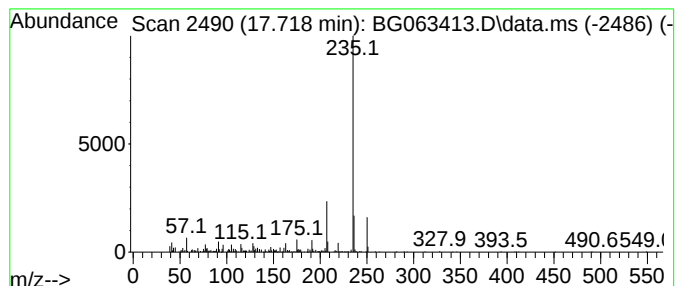
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 Benzoic acid, 3,5-bis(1,1-d... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.718	3.49 ng/ul	396480	Phenanthrene-d10	17.224

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzoic acid, 3,5-bis(1,1-dimeth...	250	C15H22O3	001421-49-4	91
2			Benzoic acid, 3,5-bis(1,1-dimeth...	250	C15H22O3	001421-49-4	87
3			2-(Ethylsulfanyl)-1H-benzimidazo...	250	C12H18N2SSi	1000486-40-4	78
4			6-Methyl-2-phenyl-4-quinolino	235	C16H13NO	001148-49-8	72
5			2,5-di-tert-Butyl-1,4-dimethoxyb...	250	C16H26O2	007323-63-9	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111324\
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Acq On : 15 Nov 2024 5:25
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Misc :
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Instrument :
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ClientSampleId :
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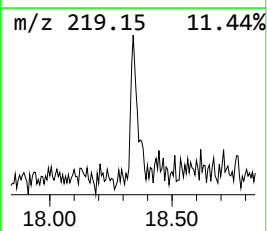
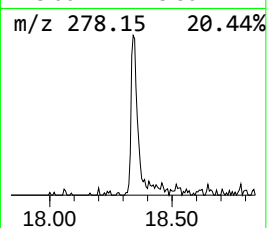
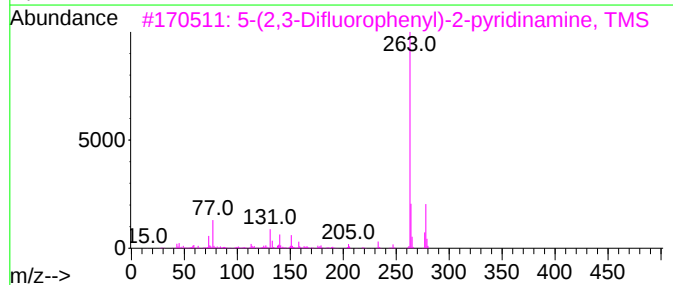
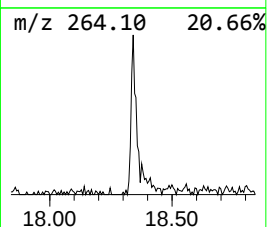
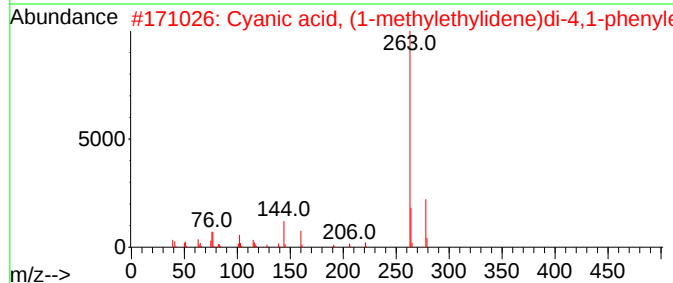
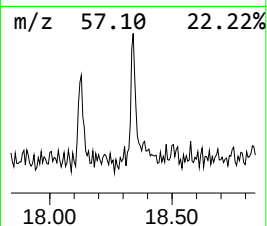
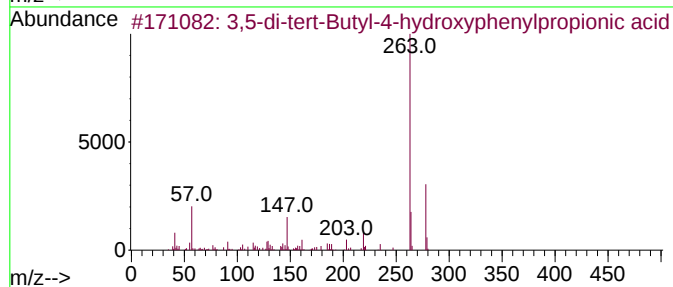
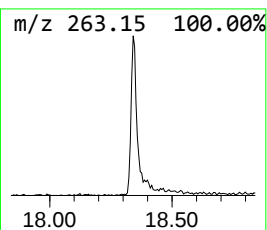
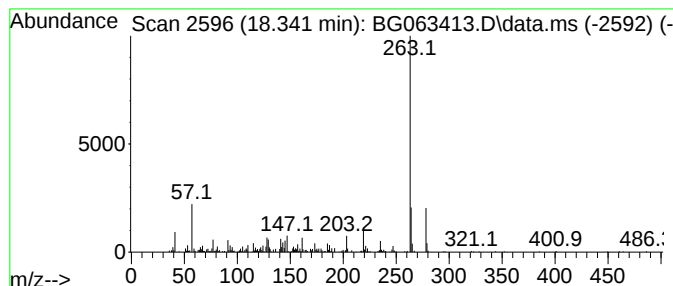
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 3,5-di-tert-Butyl-4-hydroxy... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.341	3.18 ng/ul	361233	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	86
2			Cyanic acid, (1-methylethylidene...	278	C17H14N2O2	001156-51-0	72
3			5-(2,3-Difluorophenyl)-2-pyridin...	278	C14H16F2N2Si	1000504-49-4	72
4			2,6-Bis(tert-butyl)phenol, TMS d...	278	C17H30OSi	010416-73-6	64
5			6,7-Dimethoxy-4(1H)-quinazolinon...	278	C13H18N2O3Si	1000479-65-5	56



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111324\
Data File : BG063413.D
Acq On : 15 Nov 2024 5:25
Operator : RC/JU
Sample : P4761-03
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Propanenitrile,...	7.900	25.6	ng/ul	1000300	1	7.830	779848	20.0
Triethyl phosphate	9.287	4.8	ng/ul	270510	2	10.620	1128000	20.0
4-t-Butyl-o-xylene	9.898	9.3	ng/ul	522251	2	10.620	1128000	20.0
2,4,6-Cyclohept...	10.039	3.8	ng/ul	215970	2	10.620	1128000	20.0
Benzophenone	15.832	3.9	ng/ul	375451	3	14.469	1904820	20.0
Benzenesulfonam...	15.979	2.1	ng/ul	244770	4	17.224	2272820	20.0
Benzoic acid, 3...	17.718	3.5	ng/ul	396480	4	17.224	2272820	20.0
3,5-di-tert-But...	18.341	3.2	ng/ul	361233	4	17.224	2272820	20.0