

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101724\
 Data File : BM048110.D
 Acq On : 17 Oct 2024 17:06
 Operator : RC/JU
 Sample : P4360-03
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4EB6

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_M\Methods\SFAM-EPA-BM101724.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BM048110.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.234	38	42	52	rVB	58096	93475	1.82%	0.149%
2	3.510	85	89	98	rBV	64644	109139	2.13%	0.174%
3	3.658	110	114	126	rVB	169993	308190	6.01%	0.492%
4	3.975	164	168	174	rVB3	13584	24559	0.48%	0.039%
5	4.893	321	324	332	rVB3	16628	25462	0.50%	0.041%
6	5.440	410	417	422	rBV5	11227	24387	0.48%	0.039%
7	5.857	481	488	504	rVV	512779	880189	17.16%	1.406%
8	5.987	506	510	517	rVV5	13103	32268	0.63%	0.052%
9	6.346	566	571	577	rVV5	12341	24179	0.47%	0.039%
10	6.422	577	584	595	rVV2	276481	516215	10.07%	0.825%
11	6.946	664	673	688	rVB	244492	446110	8.70%	0.713%
12	7.104	688	700	712	rBV	831054	1393542	27.17%	2.226%
13	7.304	724	734	743	rVV	899913	1520036	29.64%	2.428%
14	7.551	770	776	786	rBV	112632	201583	3.93%	0.322%
15	7.769	806	813	819	rBV	697350	1170098	22.82%	1.869%
16	7.840	819	825	835	rVV	148905	310766	6.06%	0.496%
17	7.928	835	840	846	rVB2	28605	51729	1.01%	0.083%
18	8.081	861	866	877	rVB	59025	115771	2.26%	0.185%
19	8.381	912	917	923	rBV2	32677	58738	1.15%	0.094%
20	8.481	923	934	943	rVV	720483	1269811	24.76%	2.028%
21	8.557	943	947	951	rVV	137466	238520	4.65%	0.381%
22	8.592	951	953	958	rVV	77393	102996	2.01%	0.165%
23	8.640	958	961	965	rVB2	31564	31762	0.62%	0.051%
24	8.869	993	1000	1004	rBV2	160656	298474	5.82%	0.477%
25	8.928	1004	1010	1019	rVV	895130	1562572	30.47%	2.496%
26	9.016	1019	1025	1032	rVB	370059	555873	10.84%	0.888%
27	9.098	1033	1039	1048	rBV	180053	298484	5.82%	0.477%
28	9.316	1071	1076	1080	rBV2	36266	71255	1.39%	0.114%
29	9.492	1099	1106	1114	rVB5	27319	50499	0.98%	0.081%
30	9.645	1124	1132	1138	rBV	748042	1283706	25.03%	2.050%
31	9.822	1154	1162	1168	rVV2	482001	860133	16.77%	1.374%
32	9.975	1184	1188	1192	rBV4	16617	26285	0.51%	0.042%
33	10.051	1192	1201	1210	rVV2	267612	685398	13.36%	1.095%
34	10.192	1217	1225	1244	rBV	1080890	1965321	38.32%	3.139%
35	10.345	1245	1251	1258	rVV2	61492	113558	2.21%	0.181%
36	10.416	1259	1263	1268	rVV4	32489	50997	0.99%	0.081%

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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_M\Methods\SFAM-EPA-BM101724.MA.M
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37	10.475	1268	1273	1280	rVB5	21995	49997	0.97%	0.080%
38	10.563	1280	1288	1295	rBV	935489	1780964	34.73%	2.845%
39	10.622	1295	1298	1304	rVB2	137033	202949	3.96%	0.324%
40	10.704	1304	1312	1326	rBV	246179	505037	9.85%	0.807%
41	10.851	1330	1337	1346	rBV	282139	502031	9.79%	0.802%
42	10.957	1346	1355	1365	rVB6	46443	127245	2.48%	0.203%
43	11.098	1374	1379	1385	rBV	134430	222846	4.35%	0.356%
44	11.163	1385	1390	1394	rBV	134539	213476	4.16%	0.341%
45	11.210	1394	1398	1408	rVB	66820	138800	2.71%	0.222%
46	11.533	1448	1453	1458	rBV7	16794	29213	0.57%	0.047%
47	11.704	1476	1482	1496	rBV2	114791	291957	5.69%	0.466%
48	11.839	1496	1505	1508	rVV8	30302	96720	1.89%	0.154%
49	11.881	1509	1512	1522	rVV	30968	60434	1.18%	0.097%
50	11.963	1522	1526	1535	rVB5	29100	65416	1.28%	0.104%
51	12.075	1535	1545	1547	rBV9	8610	23252	0.45%	0.037%
52	12.151	1554	1558	1563	rBV	19176	32006	0.62%	0.051%
53	12.210	1563	1568	1575	rBV	37808	77399	1.51%	0.124%
54	12.392	1591	1599	1606	rVB	23883	48742	0.95%	0.078%
55	12.539	1616	1624	1629	rBV6	15909	31174	0.61%	0.050%
56	12.769	1655	1663	1672	rBV	273918	497826	9.71%	0.795%
57	12.863	1675	1679	1690	rVV	47014	77710	1.52%	0.124%
58	12.963	1690	1696	1699	rVV2	26675	37850	0.74%	0.060%
59	13.016	1699	1705	1717	rVB	375680	552756	10.78%	0.883%
60	13.257	1742	1746	1749	rVB	21832	30386	0.59%	0.049%
61	13.316	1749	1756	1757	rBV3	17985	29677	0.58%	0.047%
62	13.333	1757	1759	1764	rVB	20848	25486	0.50%	0.041%
63	13.533	1787	1793	1803	rBV6	32835	78569	1.53%	0.125%
64	13.645	1803	1812	1821	rBV3	87078	185229	3.61%	0.296%
65	13.716	1821	1824	1832	rVB7	16219	25908	0.51%	0.041%
66	13.816	1832	1841	1847	rBV	1890107	2688344	52.42%	4.294%
67	13.945	1858	1863	1868	rVB	73993	105486	2.06%	0.168%
68	14.098	1877	1889	1897	rBV	2266033	3450039	67.27%	5.511%
69	14.274	1914	1919	1927	rVB2	57787	101802	1.99%	0.163%
70	14.410	1934	1942	1952	rBV	1315410	2018800	39.37%	3.225%
71	14.633	1974	1980	1998	rVV3	121762	342541	6.68%	0.547%
72	14.786	2000	2006	2008	rVV6	12027	24599	0.48%	0.039%
73	14.839	2008	2015	2020	rVV6	22304	54493	1.06%	0.087%
74	14.945	2028	2033	2036	rVV4	16036	27812	0.54%	0.044%
75	14.998	2038	2042	2048	rVV3	32568	47482	0.93%	0.076%
76	15.104	2052	2060	2065	rVV6	17695	41272	0.80%	0.066%

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 Title : SVOA CALIBRATION

77	15.157	2065	2069	2073	rVV	23876	35910	0.70%	0.057%
78	15.227	2073	2081	2089	rVB2	130624	219007	4.27%	0.350%
79	15.398	2103	2110	2124	rVV	2710179	4080210	79.56%	6.517%
80	15.527	2124	2132	2143	rVV	868249	1377550	26.86%	2.200%
81	15.639	2146	2151	2156	rVV6	19406	42849	0.84%	0.068%
82	15.763	2164	2172	2182	rVB	92768	148119	2.89%	0.237%
83	15.945	2198	2203	2211	rBV7	15738	31074	0.61%	0.050%
84	16.116	2225	2232	2236	rBV5	19864	42631	0.83%	0.068%
85	17.151	2401	2408	2418	rBV	1463301	2176963	42.45%	3.477%
86	17.245	2418	2424	2439	rVV	2913605	4394721	85.69%	7.020%
87	17.433	2452	2456	2463	rBV	30594	56204	1.10%	0.090%
88	17.815	2515	2521	2532	rBV	2009477	2543570	49.60%	4.063%
89	18.039	2555	2559	2567	rBV3	49100	88405	1.72%	0.141%
90	18.115	2568	2572	2573	rVV	28008	36013	0.70%	0.058%
91	18.133	2573	2575	2581	rVB4	29758	41089	0.80%	0.066%
92	18.251	2590	2595	2599	rBV	46268	73758	1.44%	0.118%
93	19.104	2735	2740	2746	rBV3	42760	65973	1.29%	0.105%
94	19.174	2748	2752	2756	rBV2	41330	62456	1.22%	0.100%
95	19.321	2773	2777	2787	rBV8	38534	85272	1.66%	0.136%
96	19.456	2796	2800	2805	rBV2	70257	95153	1.86%	0.152%
97	19.539	2807	2814	2826	rBV	3501212	5128356	100.00%	8.192%
98	21.374	3120	3126	3139	rBV	1485652	2385904	46.52%	3.811%
99	24.156	3590	3599	3615	rVB	2027926	5115959	99.76%	8.172%
100	24.356	3626	3633	3646	rVB	947152	2562452	49.97%	4.093%

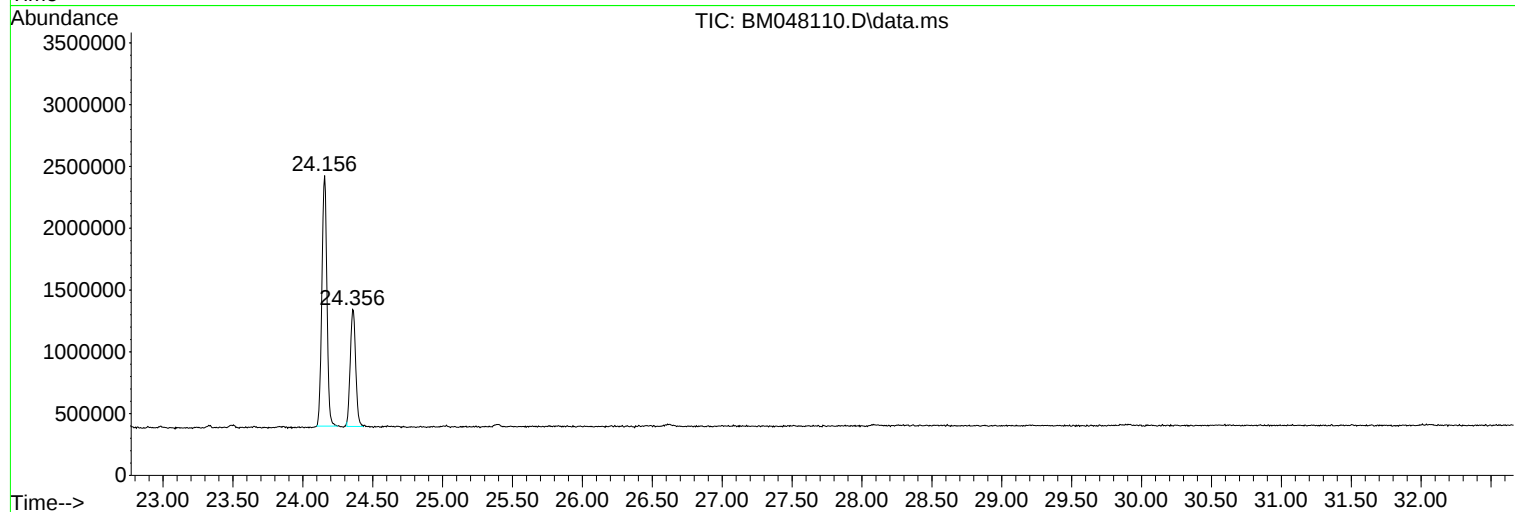
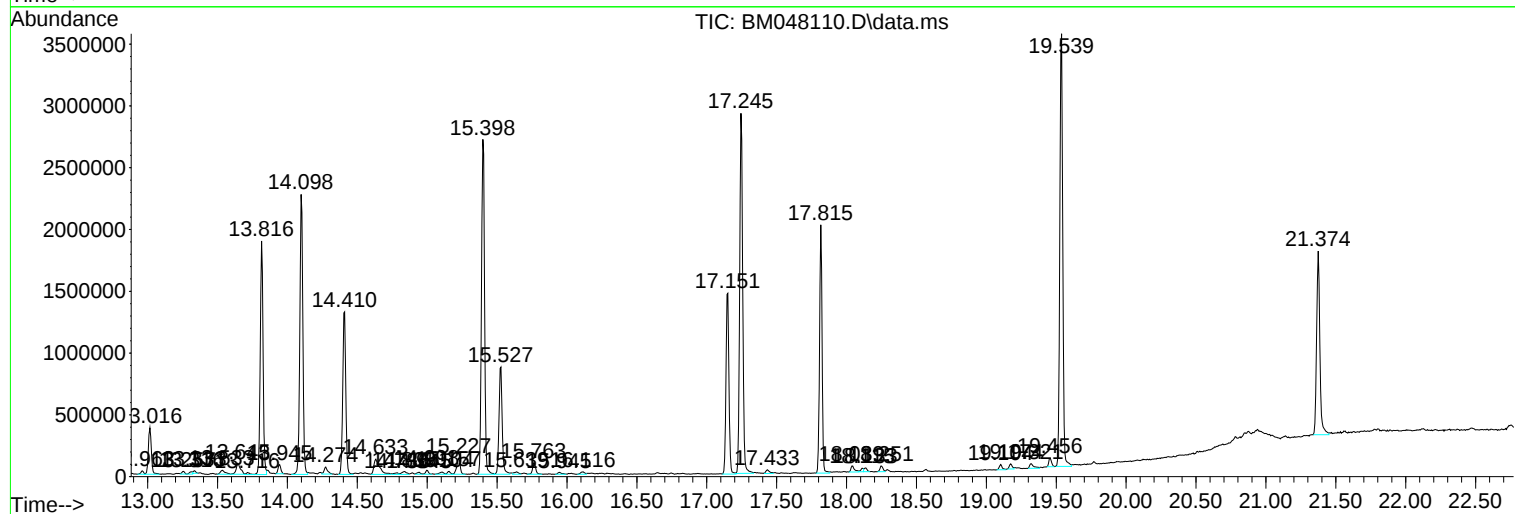
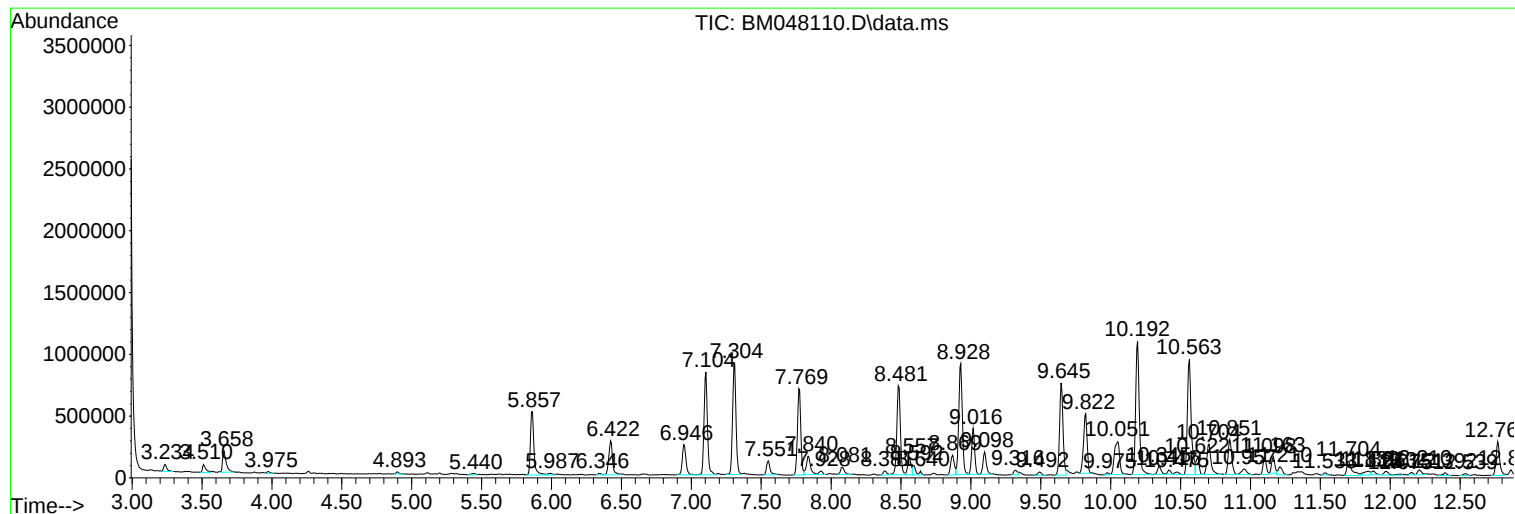
Sum of corrected areas: 62605403

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Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM101724.MA.M
Quant Title : SVOA CALIBRATION

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



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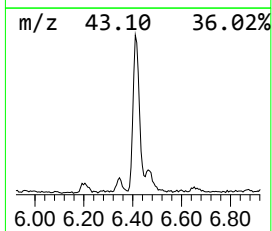
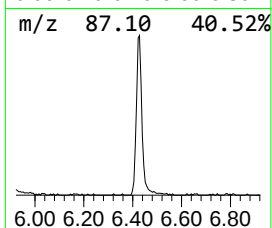
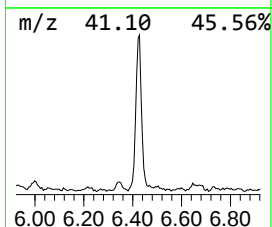
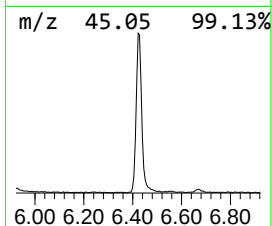
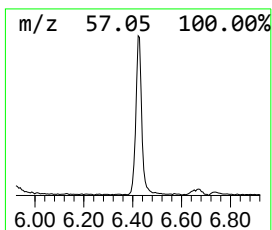
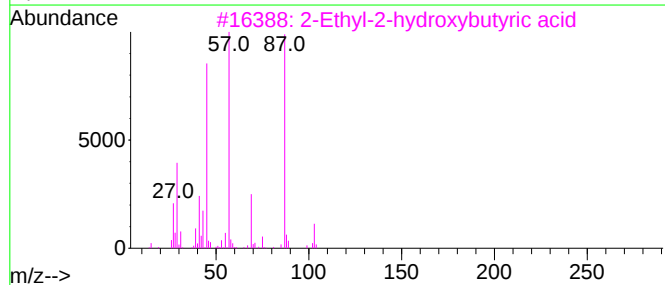
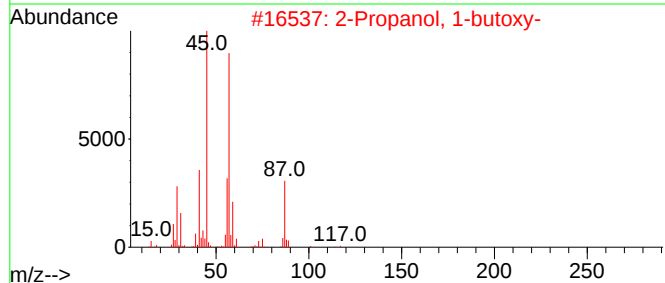
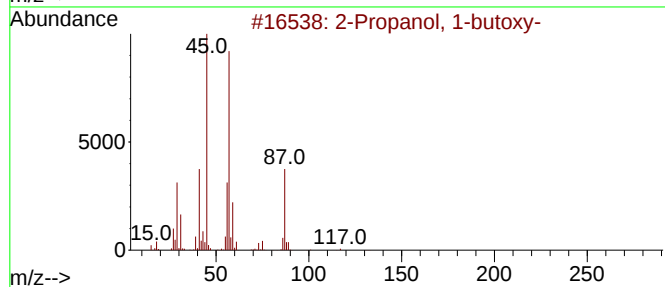
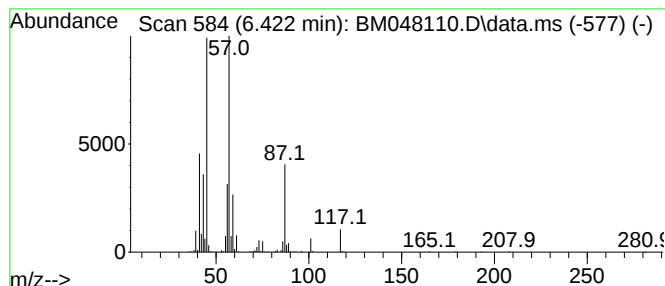
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM101724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 2-Propanol, 1-butoxy- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.422	8.82 ng/ul	516215	1,4-Dichlorobenzene-d4	7.769

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanol, 1-butoxy-			132	C7H16O2	005131-66-8	86
2	2-Propanol, 1-butoxy-			132	C7H16O2	005131-66-8	86
3	2-Ethyl-2-hydroxybutyric acid			132	C6H12O3	003639-21-2	50
4	Di-sec-Butyl ether			130	C8H18O	006863-58-7	50
5	Formic acid, 1-methylpropyl ester			102	C5H10O2	000589-40-2	50



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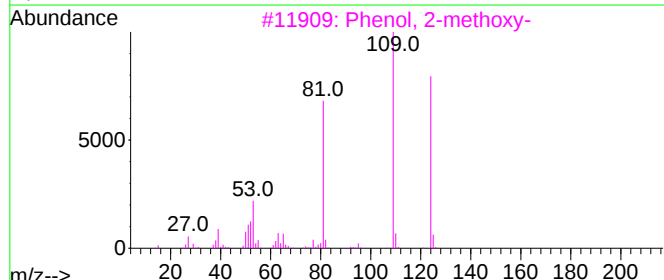
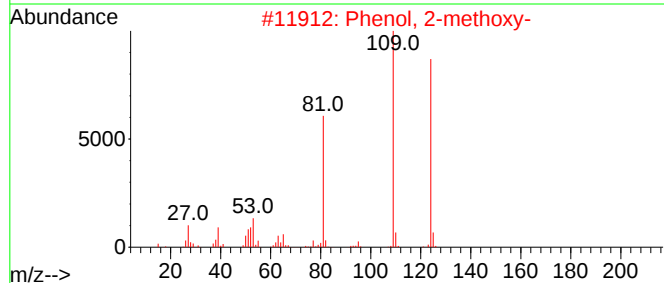
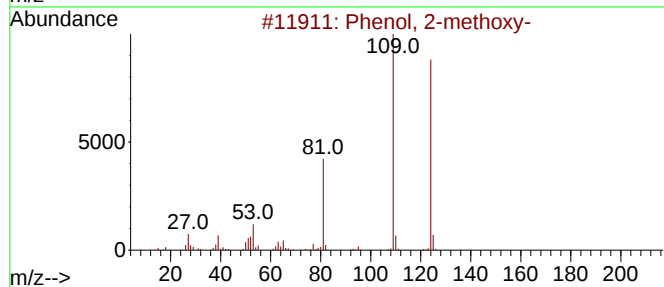
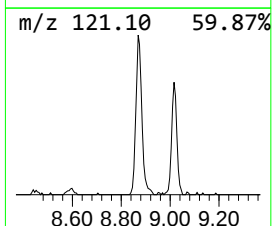
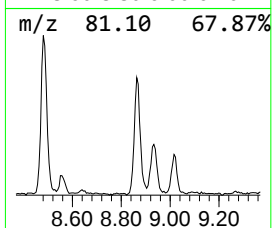
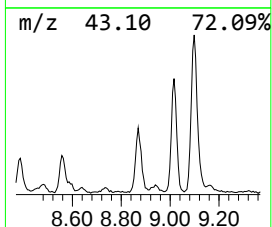
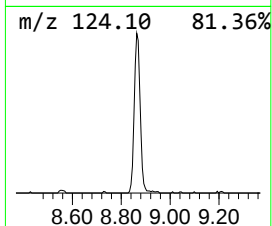
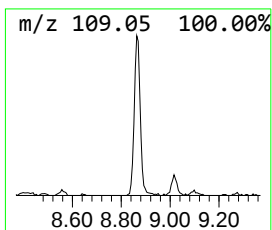
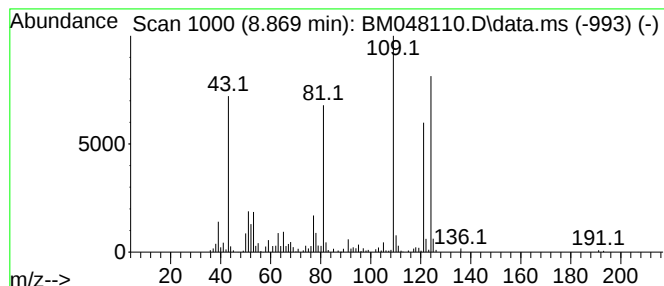
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM101724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 5 Phenol, 2-methoxy- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.869	5.10 ng/ul	298474	1,4-Dichlorobenzene-d4	7.769

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 2-methoxy-	124	C7H8O2	000090-05-1	93
2			Phenol, 2-methoxy-	124	C7H8O2	000090-05-1	93
3			Phenol, 2-methoxy-	124	C7H8O2	000090-05-1	90
4			Phenol, 2-methoxy-	124	C7H8O2	000090-05-1	87
5			Mequinol	124	C7H8O2	000150-76-5	87



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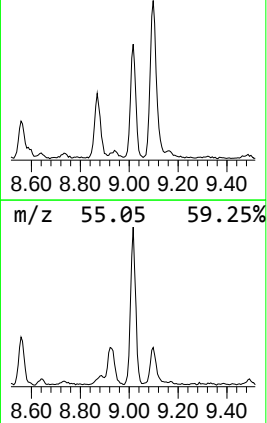
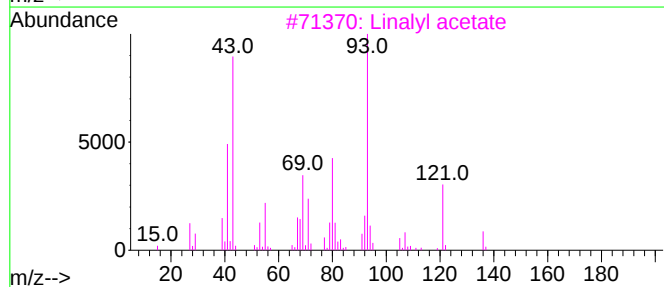
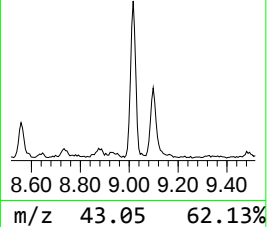
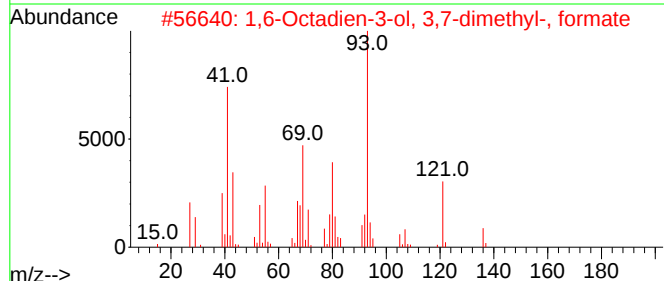
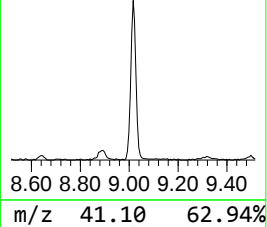
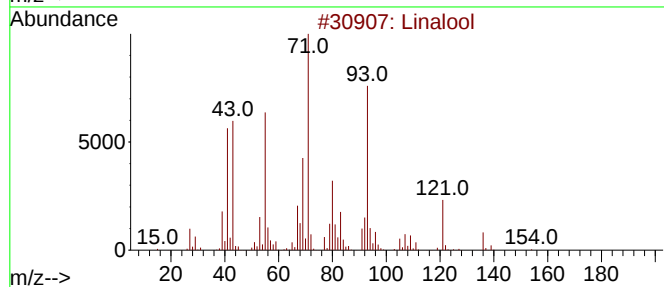
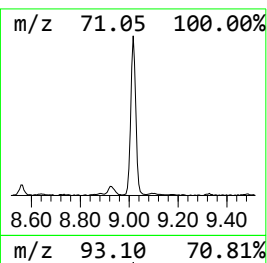
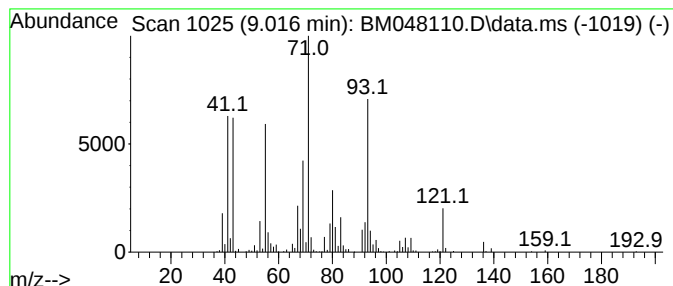
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM101724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 6 Linalool Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.016	9.50 ng/ul	555873	1,4-Dichlorobenzene-d4	7.769

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Linalool	154	C10H18O	000078-70-6	90
2			1,6-Octadien-3-ol, 3,7-dimethyl-...	182	C11H18O2	000115-99-1	49
3			Linalyl acetate	196	C12H20O2	000115-95-7	46
4			Linalyl acetate	196	C12H20O2	000115-95-7	46
5			1,3,6-Octatriene, 3,7-dimethyl-,...	136	C10H16	003338-55-4	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101724\
Data File : BM048110.D
Acq On : 17 Oct 2024 17:06
Operator : RC/JU
Sample : P4360-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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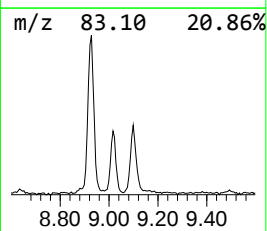
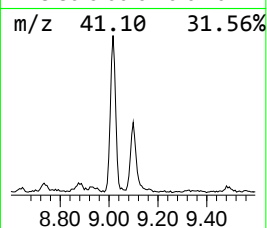
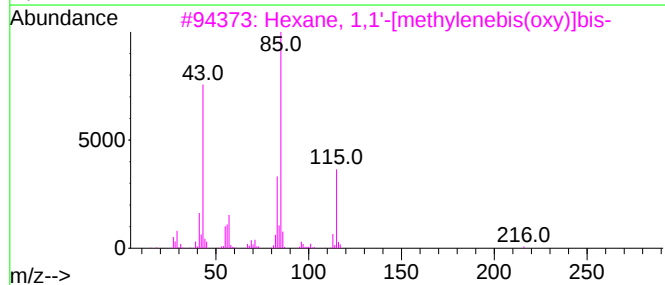
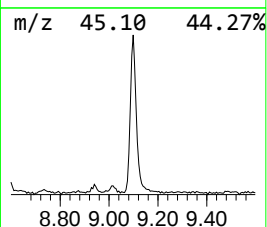
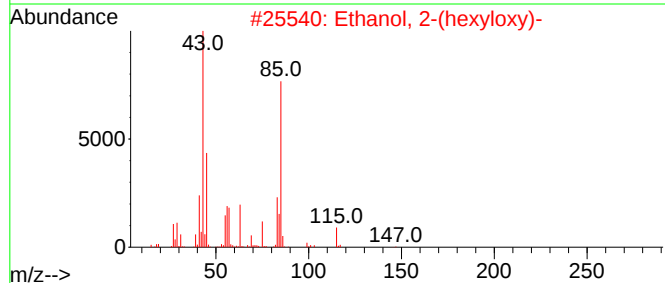
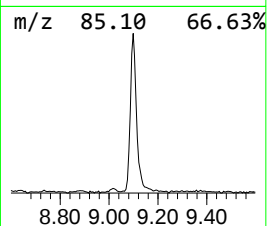
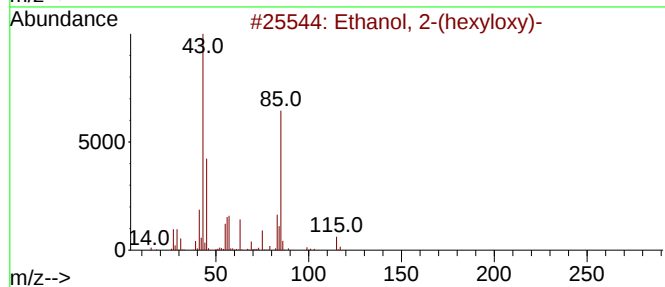
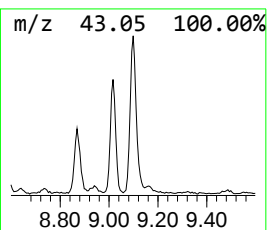
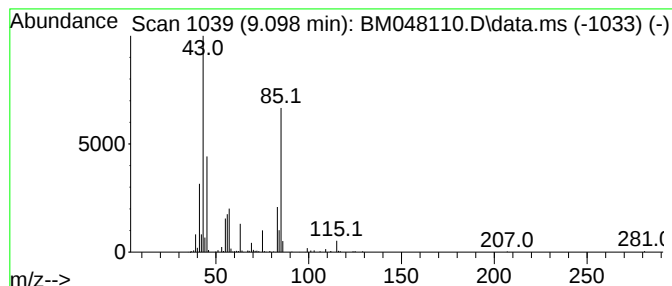
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Ethanol, 2-(hexyloxy)- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.098	5.10 ng/ul	298484	1,4-Dichlorobenzene-d4	7.769

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(hexyloxy)-	146	C8H18O2	000112-25-4	90
2			Ethanol, 2-(hexyloxy)-	146	C8H18O2	000112-25-4	59
3			Hexane, 1,1'-[methylenebis(oxy)]...	216	C13H28O2	054815-12-2	47
4			3-(Hexyloxy)-1,2-propanediol	176	C9H20O3	010305-38-1	40
5			Oxalic acid, diisohexyl ester	258	C14H26O4	1010309-32-9	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101724\
Data File : BM048110.D
Acq On : 17 Oct 2024 17:06
Operator : RC/JU
Sample : P4360-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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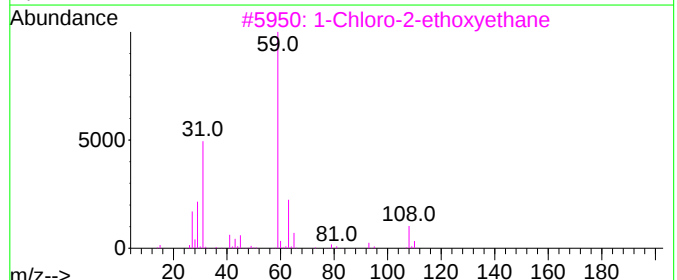
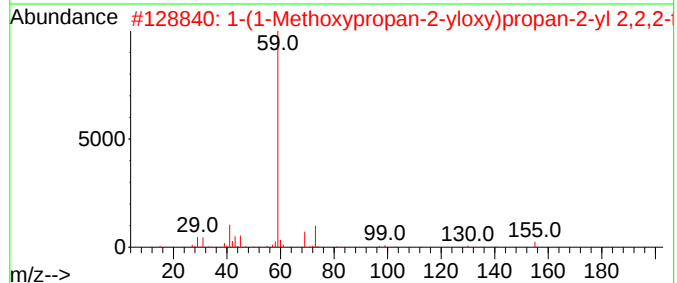
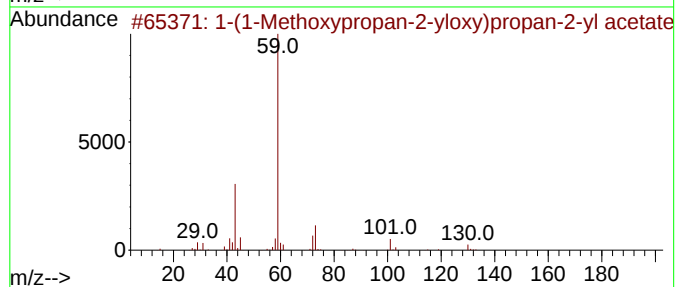
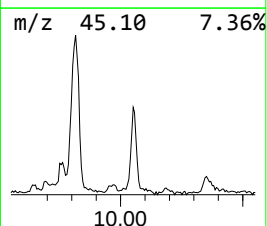
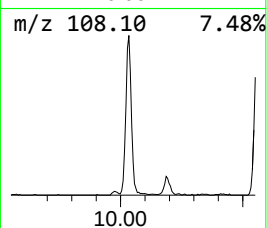
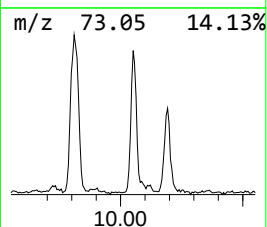
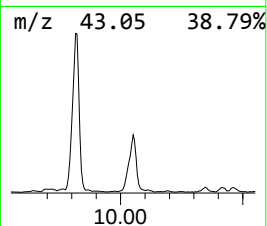
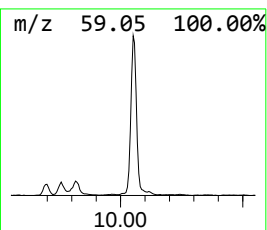
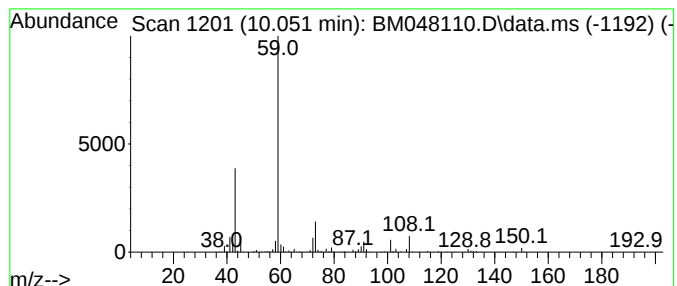
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 1-(1-Methoxypropan-2-yloxy)... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.051	7.70 ng/ul	685398	Naphthalene-d8	10.563

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-(1-Methoxypropan-2-yloxy)propa...	190	C9H18O4	1000367-08-6	86		
2	1-(1-Methoxypropan-2-yloxy)propa...	244	C9H15F3O4	1010367-08-7	64		
3	1-Chloro-2-ethoxyethane	108	C4H9ClO	000628-34-2	59		
4	Pentane, 2-methoxy-	102	C6H14O	006795-88-6	53		
5	1-(1-Methoxypropan-2-yloxy)propa...	232	C12H24O4	1010367-13-4	50		



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101724\
Data File : BM048110.D
Acq On : 17 Oct 2024 17:06
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Sample : P4360-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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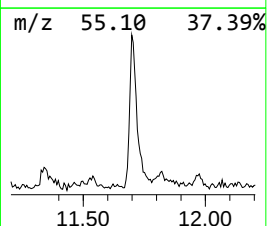
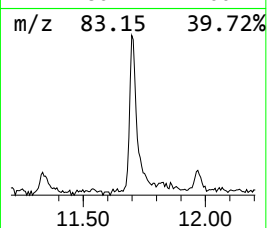
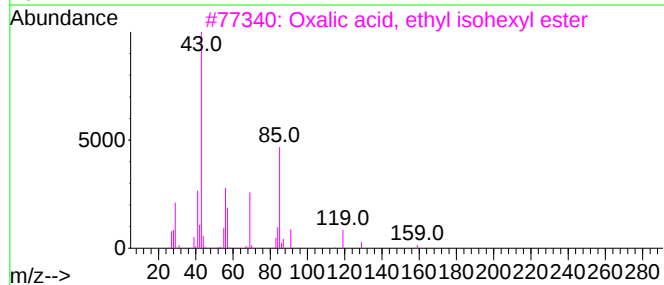
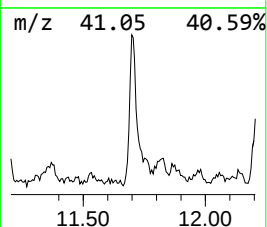
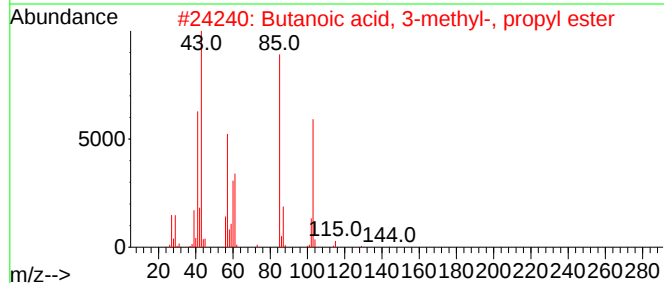
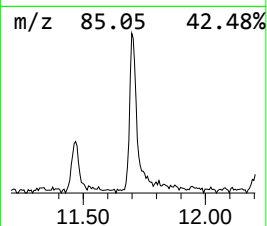
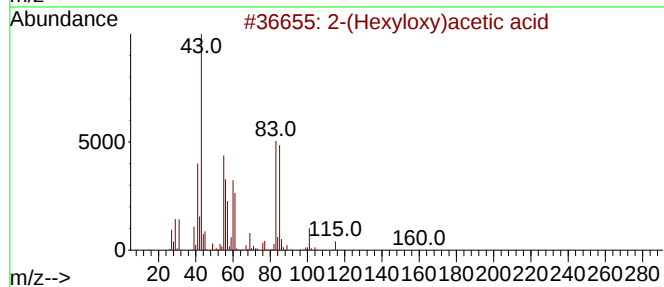
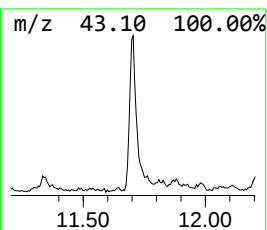
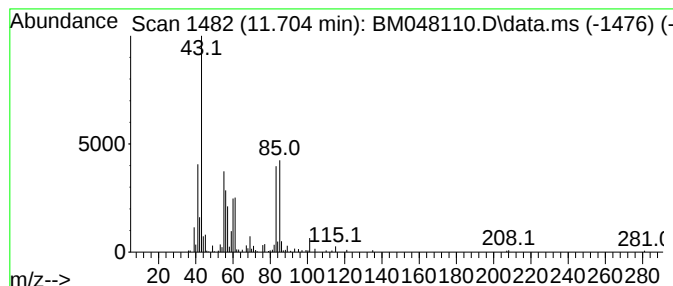
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 2-(Hexyloxy)acetic acid Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.704	3.28 ng/ul	291957	Naphthalene-d8	10.563

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-(Hexyloxy)acetic acid	160	C8H16O3	057931-25-6	90
2			Butanoic acid, 3-methyl-, propyl...	144	C8H16O2	000557-00-6	38
3			Oxalic acid, ethyl isohexyl ester	202	C10H18O4	1000309-32-5	27
4			Hexane, 2-methyl-	100	C7H16	000591-76-4	27
5			2,2-Dimethylvaleric acid	130	C7H14O2	001185-39-3	22



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101724\
Data File : BM048110.D
Acq On : 17 Oct 2024 17:06
Operator : RC/JU
Sample : P4360-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

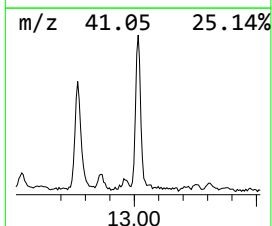
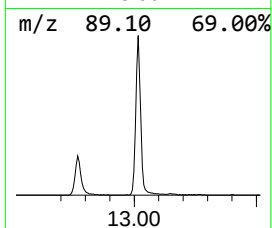
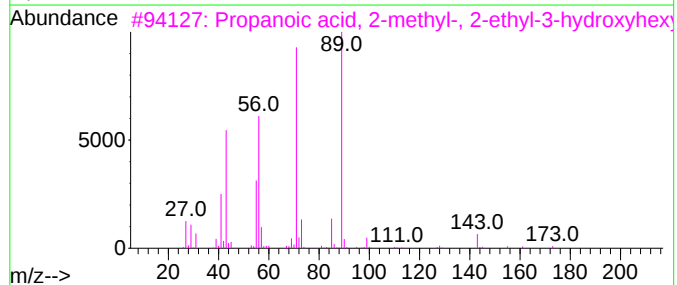
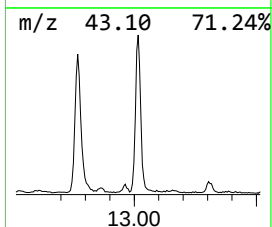
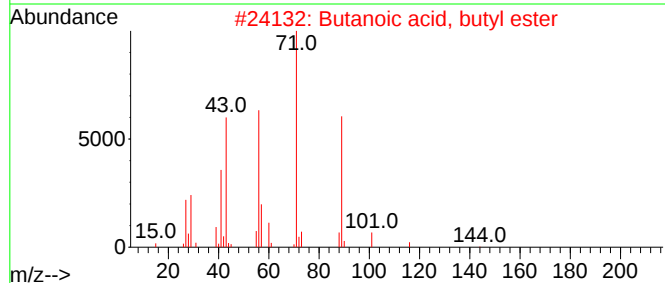
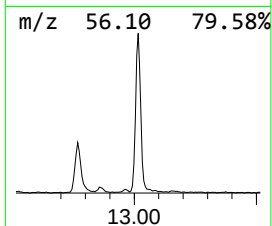
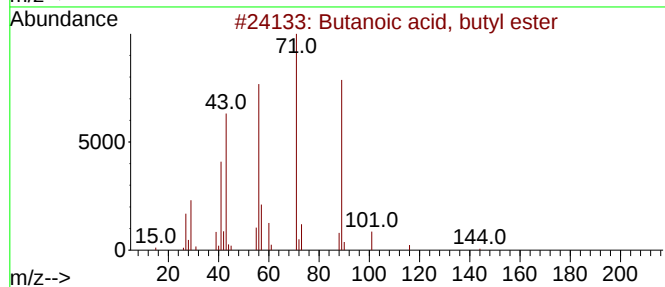
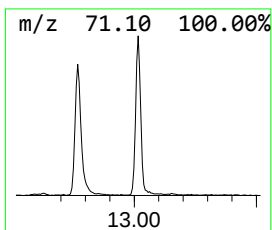
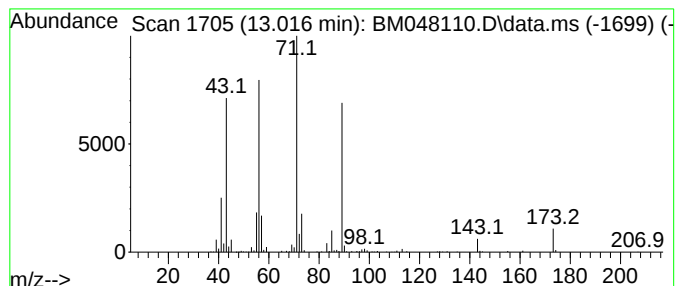
Instrument :
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 Butanoic acid, butyl ester Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.016	5.48 ng/ul	552756	Acenaphthene-d10	14.410	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butanoic acid, butyl ester	144	C8H16O2	000109-21-7	87
2	Butanoic acid, butyl ester	144	C8H16O2	000109-21-7	72
3	Propanoic acid, 2-methyl-, 2-eth...	216	C12H24O3	074367-31-0	64
4	Propanoic acid, 2-methyl-, butyl...	144	C8H16O2	000097-87-0	50
5	Propanoic acid, 2-methyl-, 2-eth...	216	C12H24O3	074367-31-0	45



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101724\
Data File : BM048110.D
Acq On : 17 Oct 2024 17:06
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Sample : P4360-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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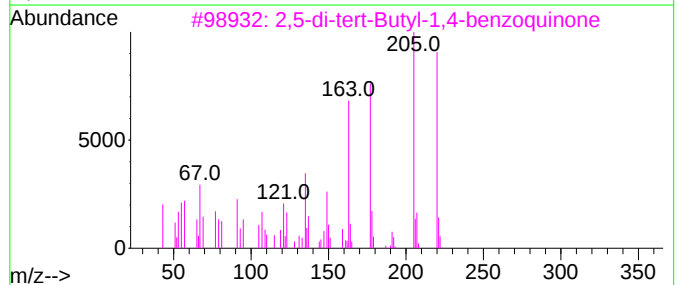
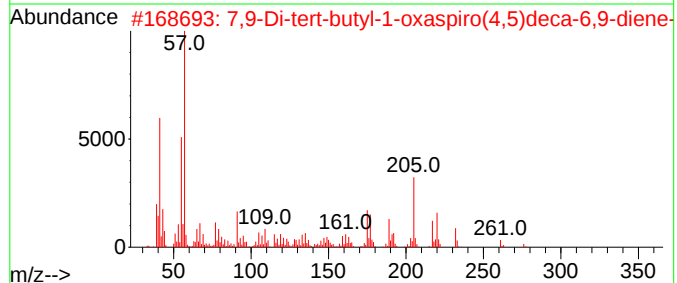
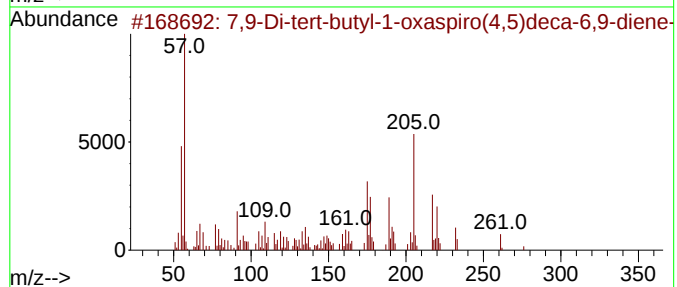
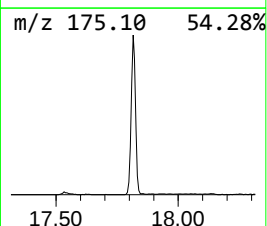
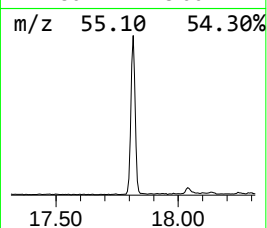
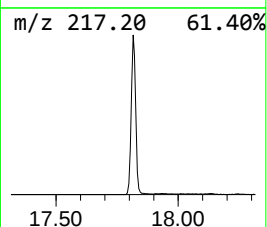
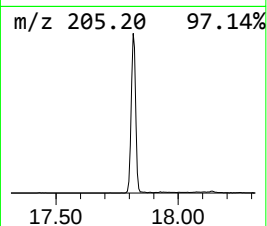
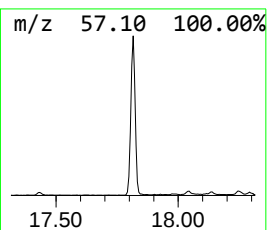
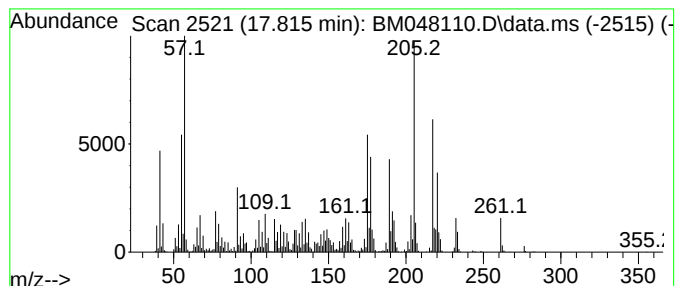
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Quant Title : SVOA CALIBRATION

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

Peak Number 17 7,9-Di-tert-butyl-1-oxaspiro... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.815	23.37 ng/ul	2543570	Phenanthrene-d10	17.151

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	7,9-Di-tert-butyl-1-oxaspiro(4,5...	276	C17H24O3	082304-66-3	99		
2	7,9-Di-tert-butyl-1-oxaspiro(4,5...	276	C17H24O3	082304-66-3	99		
3	2,5-di-tert-Butyl-1,4-benzoquinone	220	C14H20O2	002460-77-7	70		
4	7,9-Di-tert-butyl-1-oxaspiro(4,5...	276	C17H24O3	082304-66-3	55		
5	1-(3-Amino-4,6-dimethylthieno[2,...	220	C11H12N2OS	052505-42-7	50		



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101724\
Data File : BM048110.D
Acq On : 17 Oct 2024 17:06
Operator : RC/JU
Sample : P4360-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM101724.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Propanol, 1-b...	6.422	8.8	ng/ul	516215	1	7.769	1170100	20.0
Phenol, 2-methoxy-	8.869	5.1	ng/ul	298474	1	7.769	1170100	20.0
Linalool	9.016	9.5	ng/ul	555873	1	7.769	1170100	20.0
Ethanol, 2-(hex...	9.098	5.1	ng/ul	298484	1	7.769	1170100	20.0
1-(1-Methoxypro...	10.051	7.7	ng/ul	685398	2	10.563	1780960	20.0
2-(Hexyloxy)ace...	11.704	3.3	ng/ul	291957	2	10.563	1780960	20.0
Butanoic acid, ...	13.016	5.5	ng/ul	552756	3	14.410	2018800	20.0
7,9-Di-tert-but...	17.815	23.4	ng/ul	2543570	4	17.151	2176960	20.0