

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM032124\
 Data File : BM044706.D
 Acq On : 21 Mar 2024 18:41
 Operator : MA/JU
 Sample : P1813-06
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 E0LX6

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-BM030524.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BM044706.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.252	25	28	38	rVB	46105	69182	1.66%	0.150%
2	3.658	93	97	113	rBV	294902	508536	12.20%	1.101%
3	3.958	145	148	152	rVB2	12105	17541	0.42%	0.038%
4	4.393	216	222	231	rVB3	38765	62758	1.51%	0.136%
5	5.046	327	333	343	rVB2	13854	27555	0.66%	0.060%
6	5.793	454	460	463	rBV8	6484	8216	0.20%	0.018%
7	6.022	496	499	505	rVB6	4997	8851	0.21%	0.019%
8	6.199	523	529	535	rBV9	9019	17664	0.42%	0.038%
9	6.669	605	609	613	rBV5	5617	8964	0.21%	0.019%
10	6.822	629	635	637	rBV7	4874	8359	0.20%	0.018%
11	6.869	637	643	656	rVB	215039	379615	9.10%	0.822%
12	7.040	665	672	686	rBV	774642	1234297	29.60%	2.673%
13	7.175	690	695	698	rVV4	18033	33009	0.79%	0.071%
14	7.222	698	703	711	rVV	743368	1203547	28.86%	2.606%
15	7.293	713	715	719	rVB5	8903	11459	0.27%	0.025%
16	7.551	754	759	762	rBV	12125	20809	0.50%	0.045%
17	7.587	762	765	771	rVB2	9430	14288	0.34%	0.031%
18	7.693	774	783	796	rBV	637570	1051260	25.21%	2.276%
19	7.987	823	833	838	rBV4	13593	29278	0.70%	0.063%
20	8.051	838	844	851	rVV2	67119	117618	2.82%	0.255%
21	8.169	859	864	871	rBV5	13747	23990	0.58%	0.052%
22	8.322	885	890	894	rBV2	22071	39842	0.96%	0.086%
23	8.393	894	902	912	rVB	587294	1021777	24.51%	2.212%
24	8.681	945	951	959	rVB10	6747	15465	0.37%	0.033%
25	8.781	959	968	973	rBV	104590	169266	4.06%	0.366%
26	8.851	973	980	992	rVV	914763	1512322	36.27%	3.275%
27	9.016	1004	1008	1012	rBV6	10885	18273	0.44%	0.040%
28	9.069	1012	1017	1022	rVB	47234	78078	1.87%	0.169%
29	9.122	1022	1026	1030	rBV3	14168	21313	0.51%	0.046%
30	9.198	1032	1039	1046	rBV9	24972	68303	1.64%	0.148%
31	9.298	1051	1056	1063	rVB	88309	143394	3.44%	0.310%
32	9.410	1072	1075	1081	rVB6	7236	11520	0.28%	0.025%
33	9.563	1094	1101	1115	rBV	717820	1197504	28.72%	2.593%
34	9.663	1115	1118	1122	rVV5	15558	29039	0.70%	0.063%
35	9.716	1123	1127	1137	rVB	31088	63512	1.52%	0.138%
36	9.863	1146	1152	1160	rBV2	123760	233935	5.61%	0.507%

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Integrator: RTE

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Peak Location: TOP

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Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM030524.MA.M
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37	9.939	1160	1165	1170	rVB8	4110	7242	0.17%	0.016%
38	10.092	1182	1191	1199	rBV	958958	1679571	40.28%	3.637%
39	10.169	1202	1204	1210	rVB3	15782	22387	0.54%	0.048%
40	10.298	1222	1226	1229	rVB5	5119	7655	0.18%	0.017%
41	10.469	1245	1255	1269	rVB	876016	1530998	36.72%	3.315%
42	10.622	1274	1281	1297	rVB	527151	1022946	24.53%	2.215%
43	10.828	1314	1316	1321	rVB7	7464	8959	0.21%	0.019%
44	11.016	1342	1348	1353	rBV9	8384	19495	0.47%	0.042%
45	11.175	1370	1375	1382	rVB4	19143	39368	0.94%	0.085%
46	11.245	1382	1387	1389	rBV6	6950	9183	0.22%	0.020%
47	11.375	1405	1409	1416	rVB3	14461	27863	0.67%	0.060%
48	11.545	1433	1438	1442	rBV4	10096	18583	0.45%	0.040%
49	11.657	1453	1457	1464	rBV10	7715	16007	0.38%	0.035%
50	11.822	1474	1485	1494	rBV2	58014	163282	3.92%	0.354%
51	12.069	1516	1527	1536	rVV6	20022	74398	1.78%	0.161%
52	12.245	1554	1557	1560	rBV4	5546	7368	0.18%	0.016%
53	12.286	1560	1564	1568	rVB7	6671	10579	0.25%	0.023%
54	12.375	1575	1579	1582	rBV6	5344	9308	0.22%	0.020%
55	12.416	1582	1586	1590	rVB3	16645	27115	0.65%	0.059%
56	12.510	1595	1602	1604	rBV8	10315	21651	0.52%	0.047%
57	12.610	1618	1619	1626	rVB7	8452	13187	0.32%	0.029%
58	12.704	1632	1635	1637	rBV2	8122	12055	0.29%	0.026%
59	12.798	1646	1651	1654	rBV5	10958	17951	0.43%	0.039%
60	12.828	1654	1656	1662	rVV6	11109	16639	0.40%	0.036%
61	12.933	1668	1674	1681	rVB10	14160	30523	0.73%	0.066%
62	13.039	1688	1692	1697	rVB6	9215	12778	0.31%	0.028%
63	13.116	1701	1705	1706	rBV4	11812	14873	0.36%	0.032%
64	13.175	1711	1715	1719	rVV	27976	43714	1.05%	0.095%
65	13.227	1722	1724	1733	rVB5	15470	27696	0.66%	0.060%
66	13.304	1733	1737	1744	rBV8	11488	20172	0.48%	0.044%
67	13.386	1747	1751	1757	rVB9	11294	21364	0.51%	0.046%
68	13.457	1758	1763	1764	rBV5	5849	9275	0.22%	0.020%
69	13.516	1769	1773	1777	rVV7	7581	12055	0.29%	0.026%
70	13.557	1777	1780	1784	rBV5	13409	17985	0.43%	0.039%
71	13.651	1792	1796	1798	rBV5	6527	8181	0.20%	0.018%
72	13.757	1808	1814	1821	rBV	1752167	2408902	57.77%	5.216%
73	13.957	1843	1848	1850	rBV6	13330	18128	0.43%	0.039%
74	14.027	1850	1860	1871	rVB	2052658	3086396	74.02%	6.683%
75	14.139	1876	1879	1882	rBV2	9899	11635	0.28%	0.025%
76	14.216	1887	1892	1897	rBV3	38907	68747	1.65%	0.149%

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 Title : SVOA CALIBRATION

77	14.333	1905	1912	1920	rBV	1116646	1713539	41.10%	3.710%
78	14.563	1946	1951	1961	rBV2	106717	208881	5.01%	0.452%
79	14.722	1973	1978	1984	rBV8	24572	59758	1.43%	0.129%
80	14.874	2002	2004	2010	rVB7	9743	14115	0.34%	0.031%
81	15.333	2075	2082	2095	rBV	2485206	3826167	91.76%	8.284%
82	15.463	2098	2104	2110	rBV	868875	1183599	28.39%	2.563%
83	15.580	2119	2124	2134	rVB3	100771	216410	5.19%	0.469%
84	15.863	2166	2172	2181	rBV3	69372	167957	4.03%	0.364%
85	16.045	2199	2203	2209	rBV2	20047	31801	0.76%	0.069%
86	16.110	2211	2214	2220	rVB7	21676	33966	0.81%	0.074%
87	16.386	2256	2261	2263	rBV	39107	72734	1.74%	0.157%
88	16.504	2277	2281	2287	rVB8	20367	35994	0.86%	0.078%
89	17.027	2366	2370	2373	rBV5	22416	30117	0.72%	0.065%
90	17.092	2374	2381	2388	rVV	1284373	1813527	43.49%	3.927%
91	17.192	2391	2398	2408	rBV	2535021	3677290	88.19%	7.962%
92	17.686	2479	2482	2488	rBV6	24756	40939	0.98%	0.089%
93	18.004	2531	2536	2545	rBV	151303	263240	6.31%	0.570%
94	18.445	2608	2611	2615	rVB4	47409	54508	1.31%	0.118%
95	19.292	2752	2755	2764	rVB2	70621	94786	2.27%	0.205%
96	19.492	2783	2789	2796	rBV	3081767	4169605	100.00%	9.028%
97	19.551	2796	2799	2805	rVB	172467	231023	5.54%	0.500%
98	21.292	3090	3095	3102	rBV	1421583	1941535	46.56%	4.204%
99	23.392	3446	3452	3460	rVB2	2234036	4126409	98.96%	8.935%
100	23.533	3470	3476	3488	rVB	1086506	2100225	50.37%	4.547%

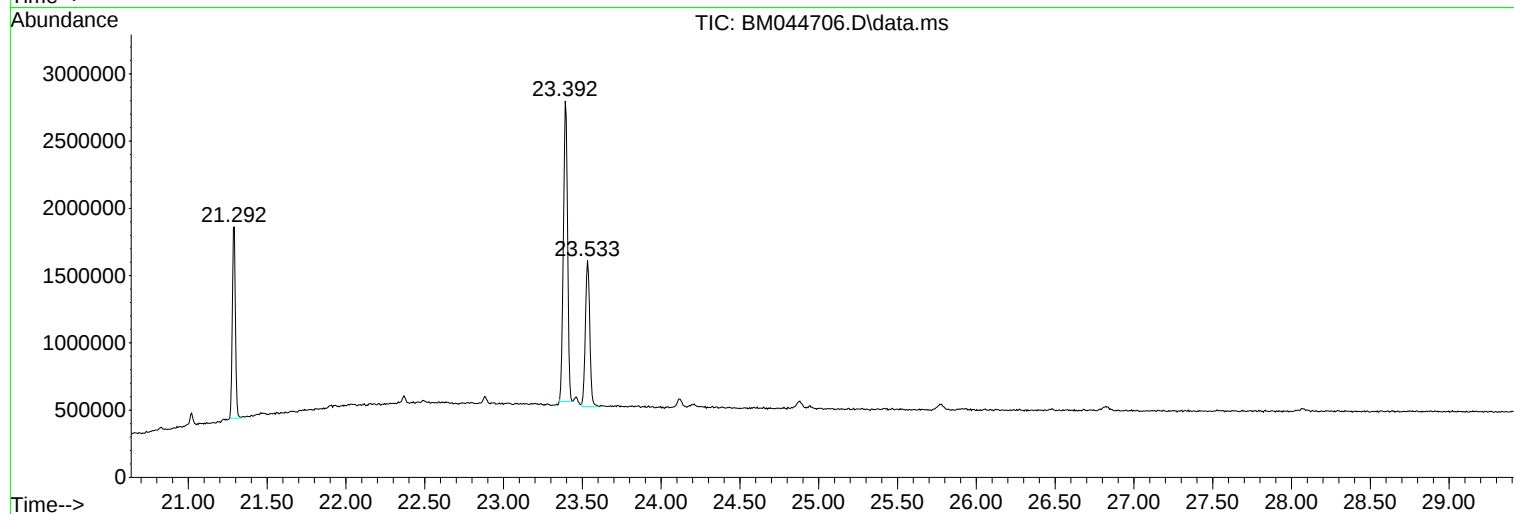
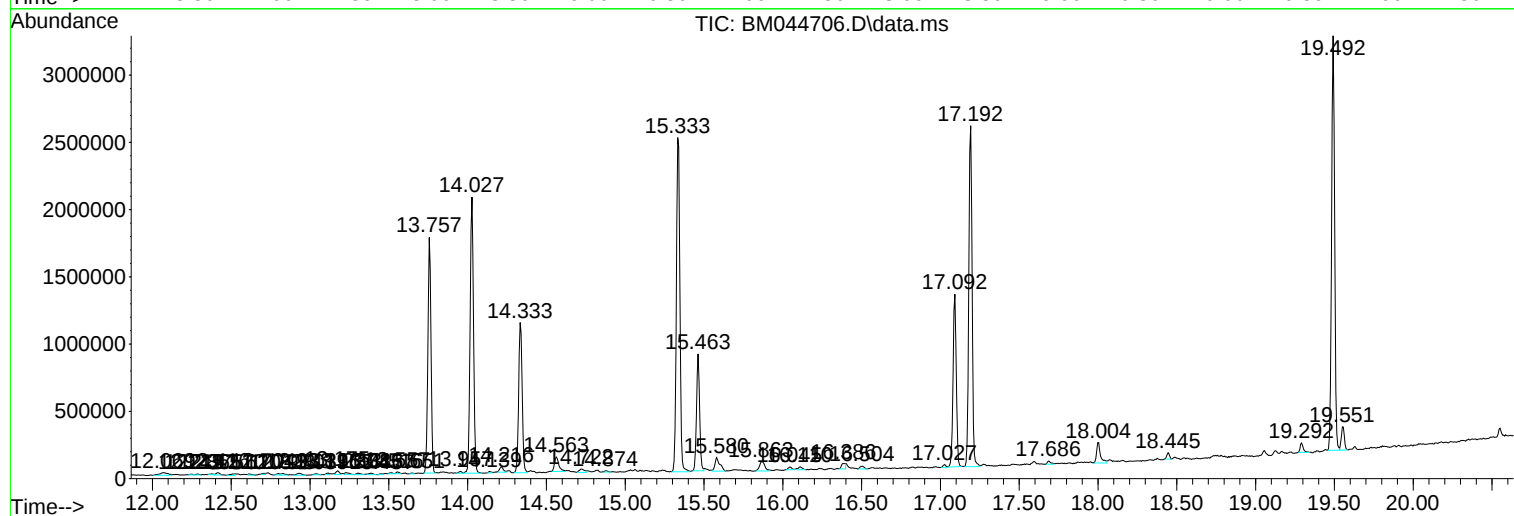
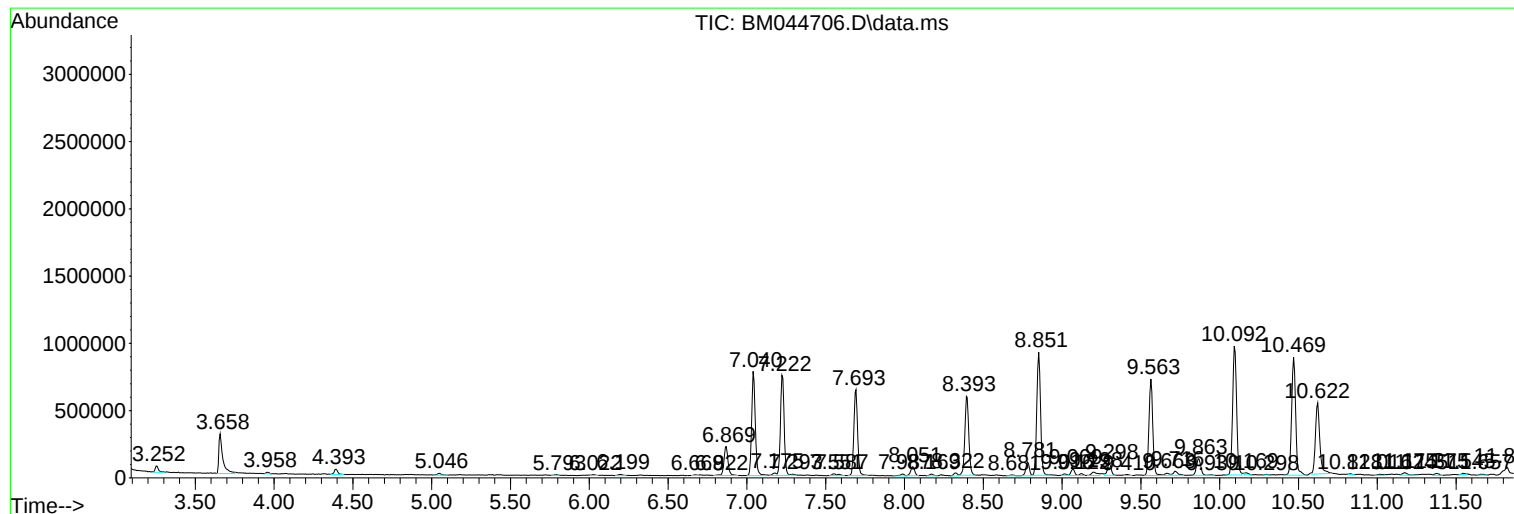
Sum of corrected areas: 46184678

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM030524.MA.M
Quant Title : SVOA CALIBRATION

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



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Instrument :
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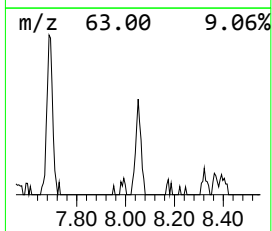
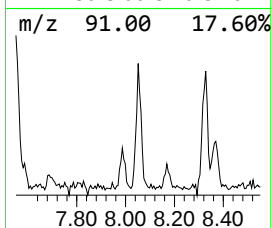
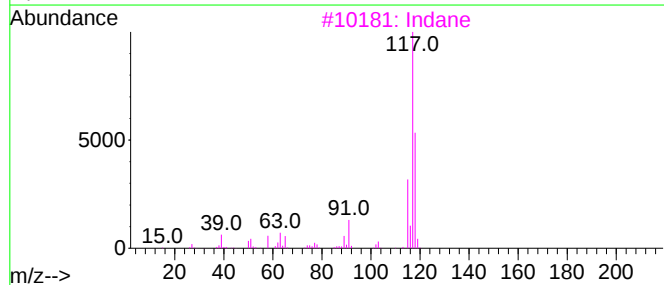
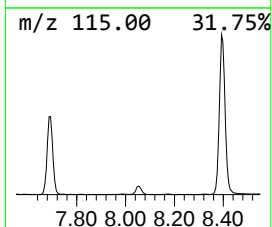
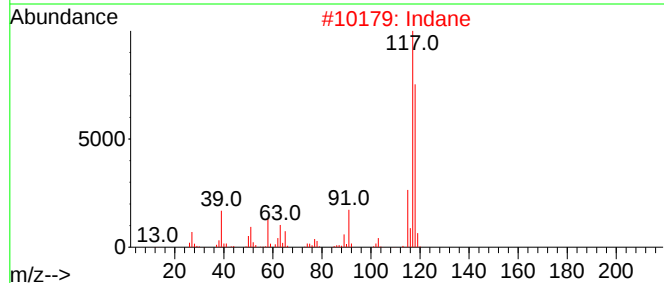
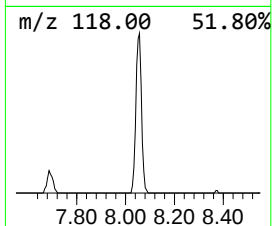
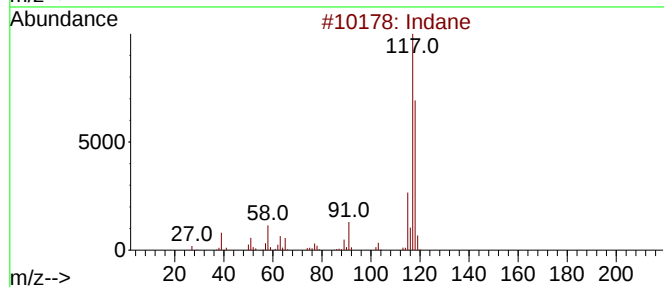
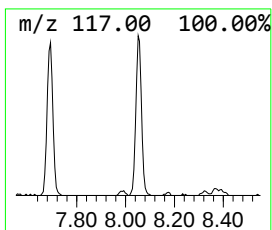
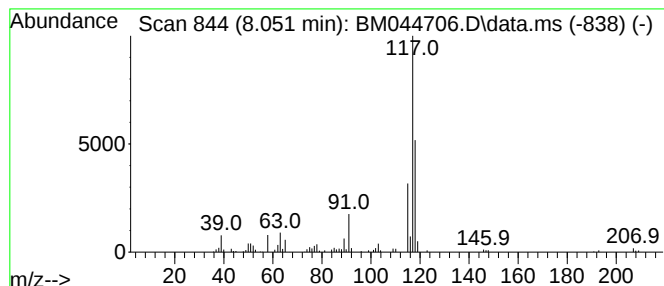
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM030524.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Indane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.051	2.24 ng/ul	117618	1,4-Dichlorobenzene-d4	7.693

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane			118	C9H10	000496-11-7	93
2	Indane			118	C9H10	000496-11-7	91
3	Indane			118	C9H10	000496-11-7	87
4	Benzene, 1-ethenyl-2-methyl-			118	C9H10	000611-15-4	81
5	Benzene, cyclopropyl-			118	C9H10	000873-49-4	74



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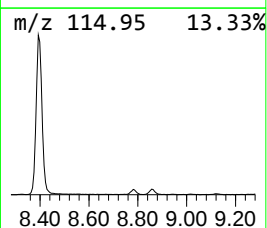
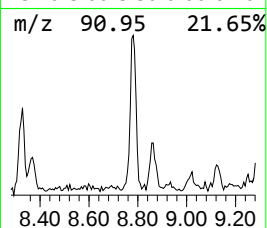
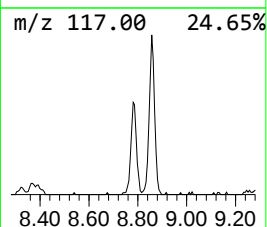
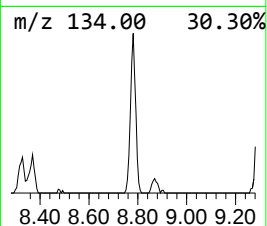
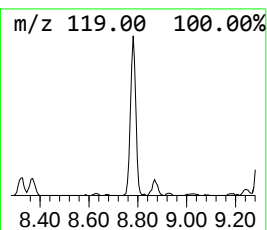
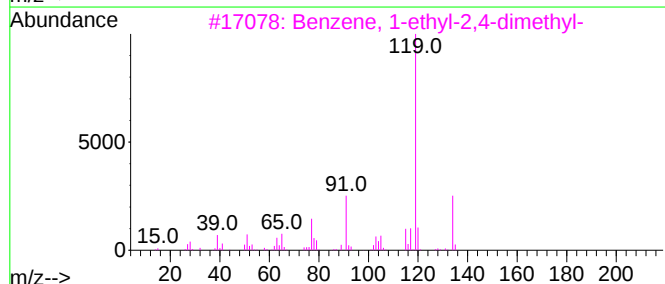
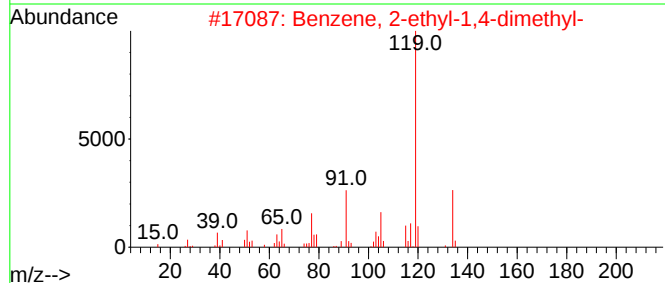
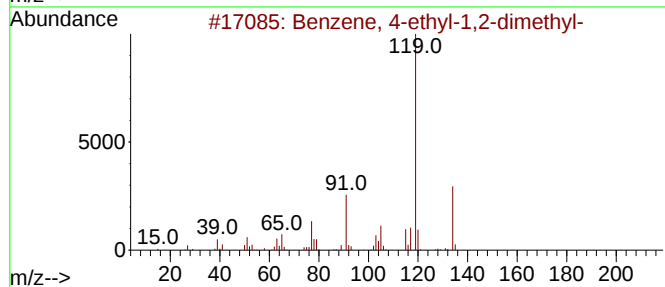
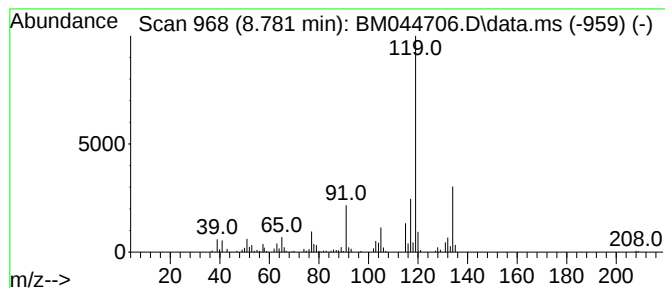
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM030524.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.781	3.22 ng/ul	169266	1,4-Dichlorobenzene-d4	7.693

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	93
2			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	90
3			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	81
4			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	81
5			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	81



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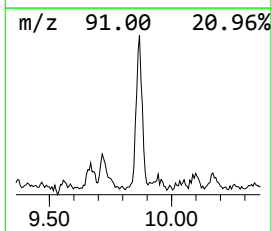
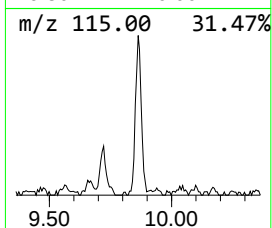
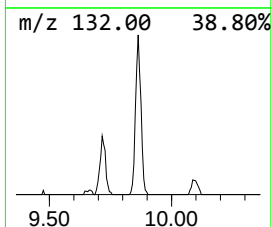
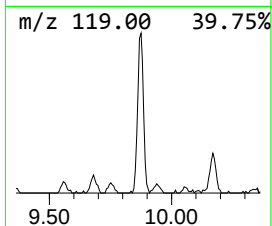
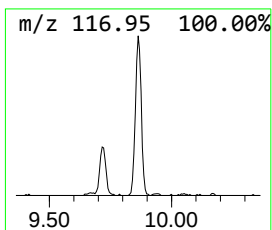
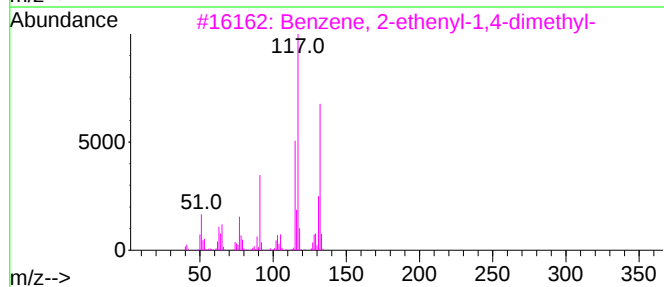
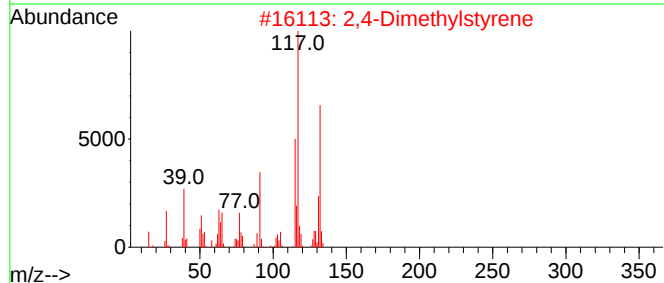
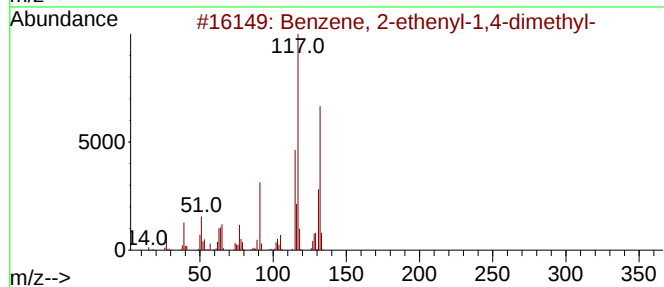
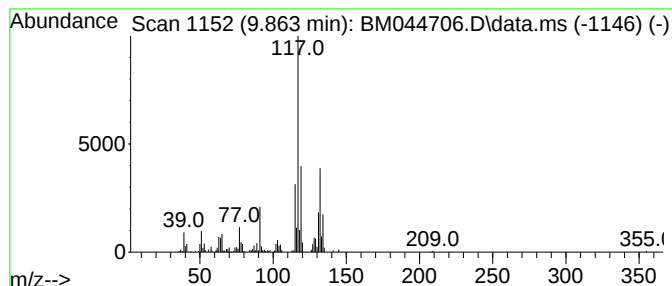
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM030524.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.863	3.06 ng/ul	233935	Naphthalene-d8	10.469

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	95
2			2,4-Dimethylstyrene	132	C10H12	002234-20-0	95
3			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	95
4			1-Phenyl-1-butene	132	C10H12	000824-90-8	89
5			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	89



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM032124\
Data File : BM044706.D
Acq On : 21 Mar 2024 18:41
Operator : MA/JU
Sample : P1813-06
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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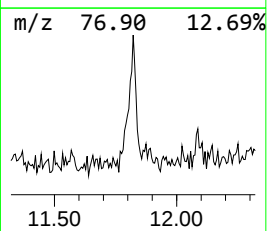
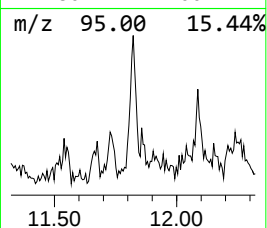
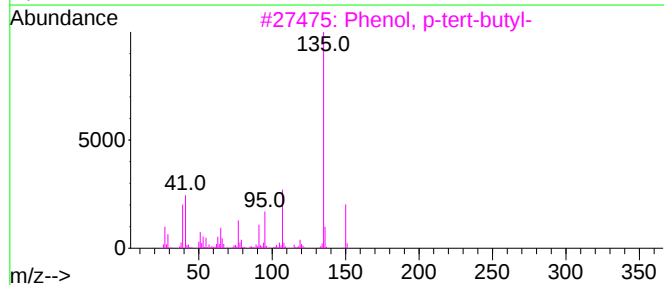
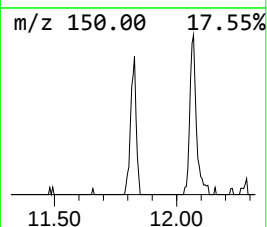
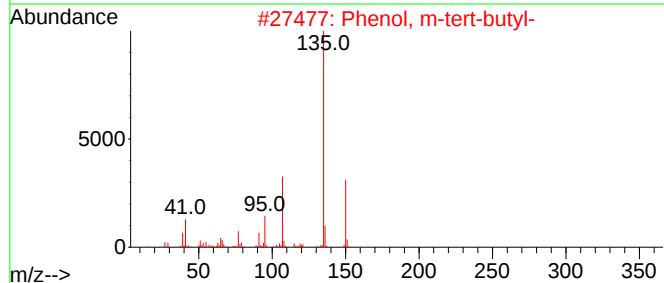
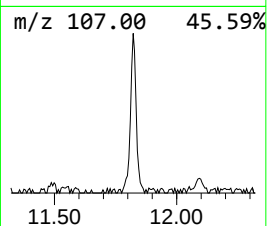
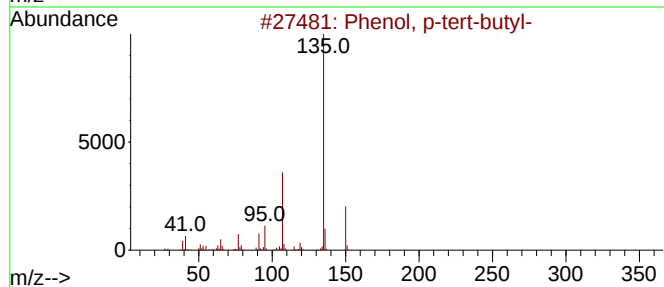
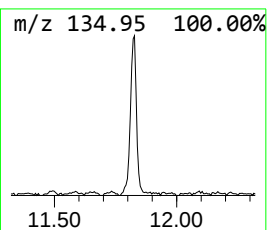
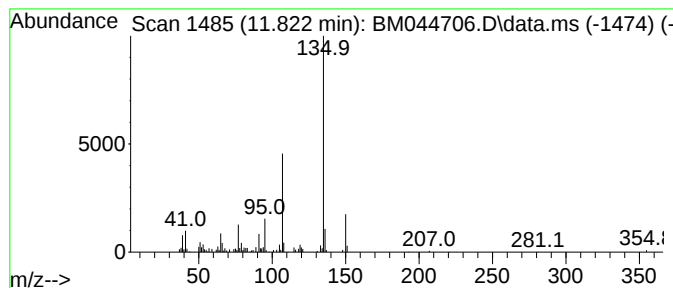
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Phenol, p-tert-butyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.822	2.13 ng/ul	163282	Naphthalene-d8	10.469

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	95
2			Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	94
3			Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	94
4			Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	91
5			Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM032124\
Data File : BM044706.D
Acq On : 21 Mar 2024 18:41
Operator : MA/JU
Sample : P1813-06
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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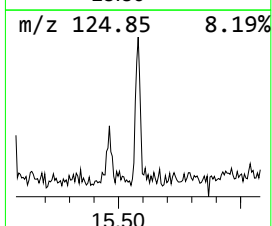
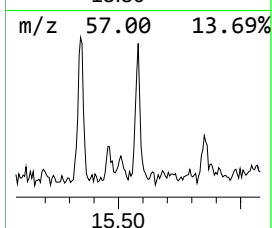
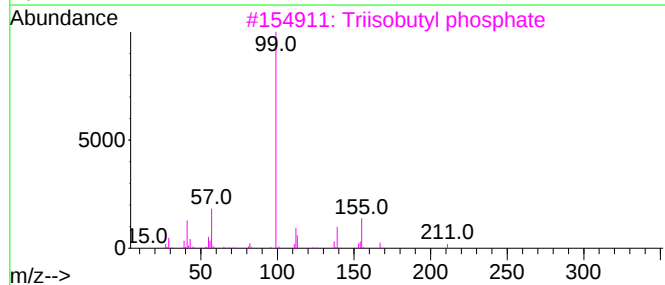
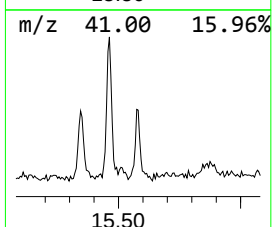
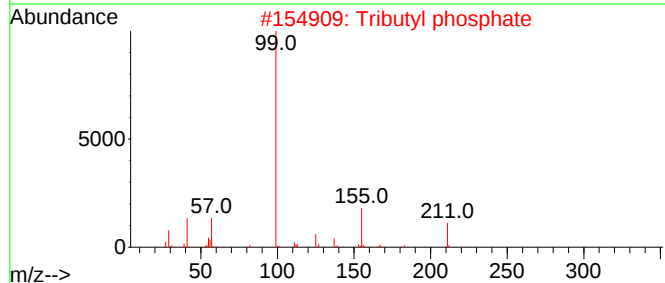
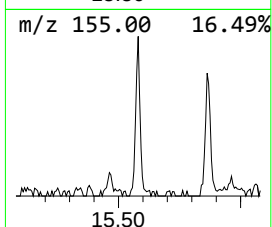
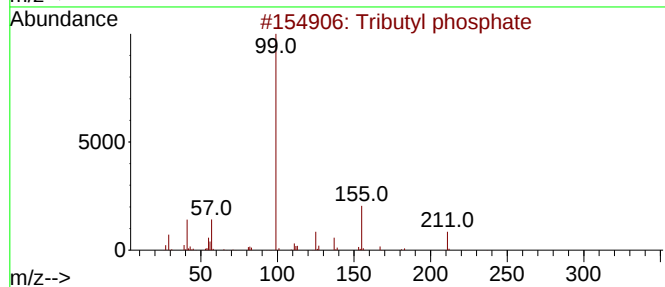
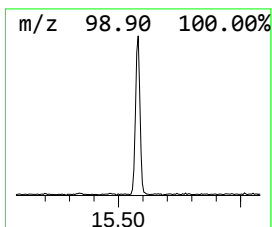
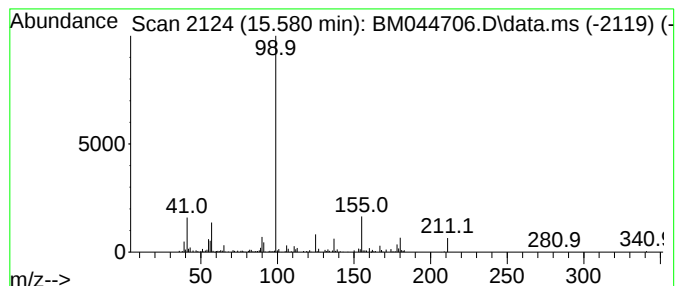
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Tributyl phosphate Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.580	2.53 ng/ul	216410	Acenaphthene-d10	14.333

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tributyl phosphate	266	C12H27O4P	000126-73-8	83
2			Tributyl phosphate	266	C12H27O4P	000126-73-8	64
3			Triisobutyl phosphate	266	C12H27O4P	000126-71-6	50
4			Phosphoric acid, dibutyl 1,1-dim...	352	C13H25F4O4P	1000298-93-9	50
5			Tributyl phosphate	266	C12H27O4P	000126-73-8	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM032124\
Data File : BM044706.D
Acq On : 21 Mar 2024 18:41
Operator : MA/JU
Sample : P1813-06
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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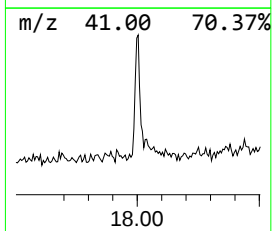
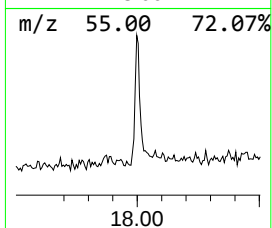
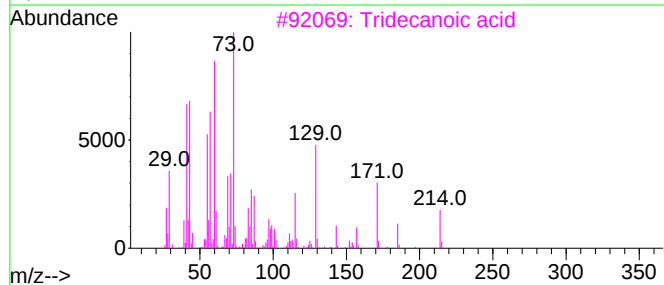
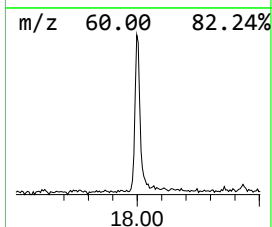
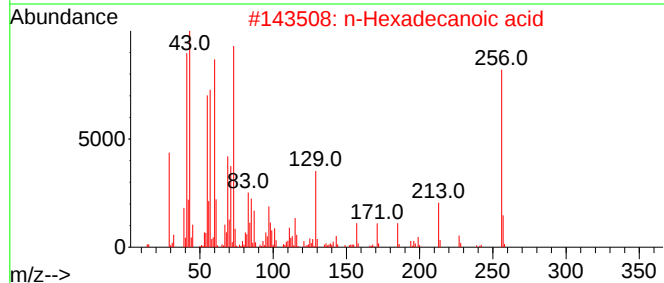
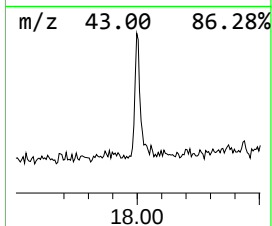
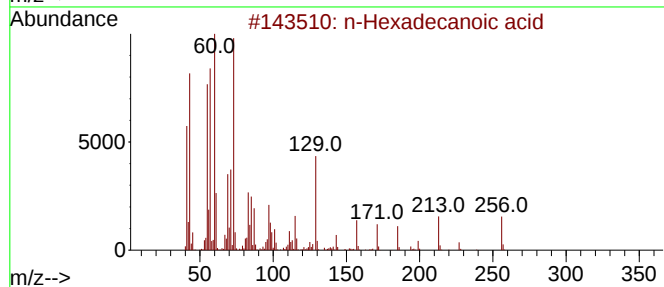
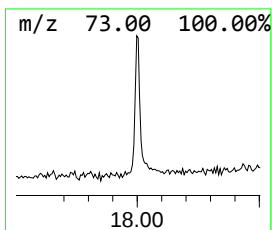
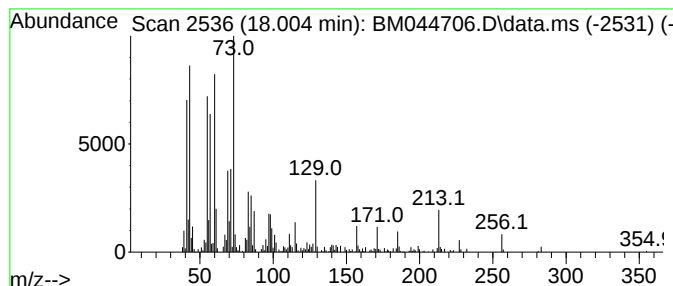
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.004	2.90 ng/ul	263240	Phenanthrene-d10	17.092

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
3			Tridecanoic acid	214	C13H26O2	000638-53-9	91
4			Pentadecanoic acid	242	C15H30O2	001002-84-2	87
5			Tetradecanoic acid	228	C14H28O2	000544-63-8	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM032124\
Data File : BM044706.D
Acq On : 21 Mar 2024 18:41
Operator : MA/JU
Sample : P1813-06
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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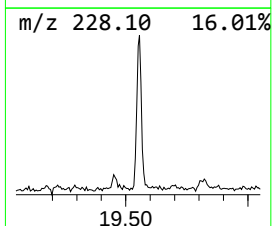
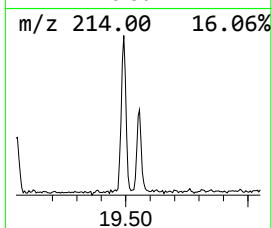
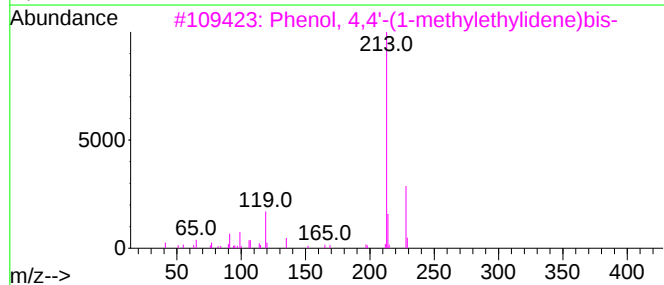
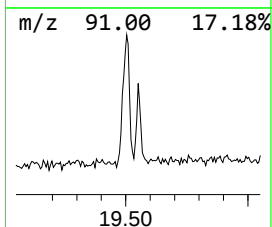
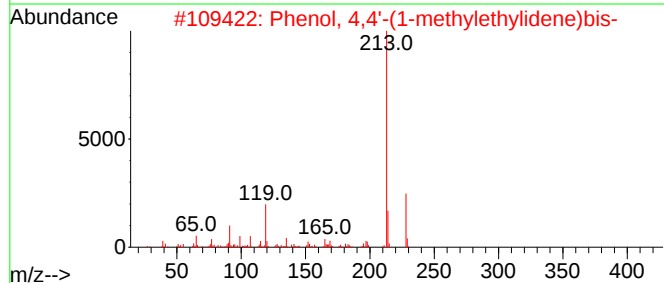
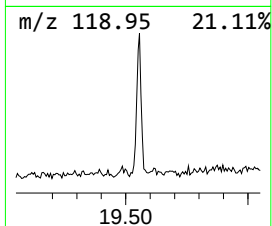
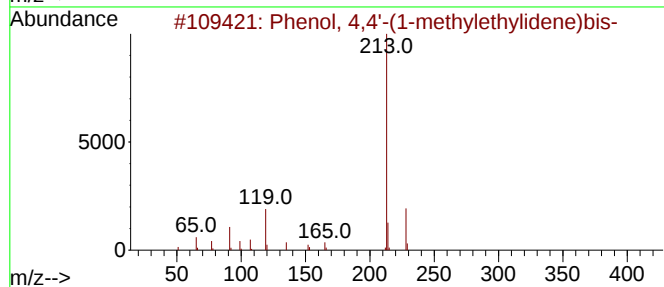
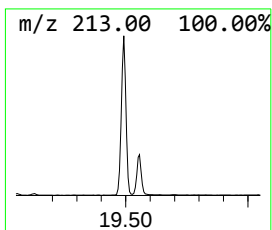
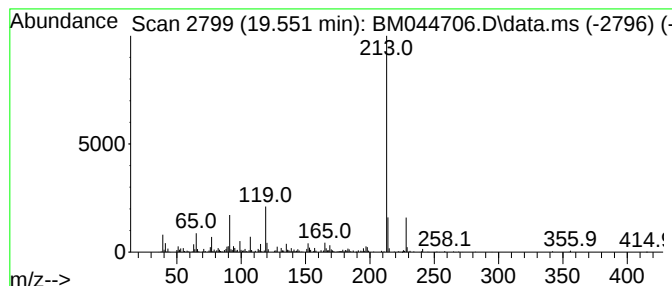
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Phenol, 4,4'-(1-methylethyl... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.551	2.38 ng/ul	231023	Chrysene-d12	21.292

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 4,4'-(1-methylethylidene...	228	C15H16O2	000080-05-7	95
2			Phenol, 4,4'-(1-methylethylidene...	228	C15H16O2	000080-05-7	90
3			Phenol, 4,4'-(1-methylethylidene...	228	C15H16O2	000080-05-7	90
4			Phenol, 4,4'-(1-methylethylidene...	228	C15H16O2	000080-05-7	87
5			4,4'-(1,3-Dimethylbutylidene)bis...	270	C18H22O2	006807-17-6	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM032124\
Data File : BM044706.D
Acq On : 21 Mar 2024 18:41
Operator : MA/JU
Sample : P1813-06
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM030524.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
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Benzene, 4-ethy...	8.781	3.2	ng/ul	169266	1	7.693	1051260	20.0
Benzene, 2-ethe...	9.863	3.1	ng/ul	233935	2	10.469	1531000	20.0
Phenol, p-tert-...	11.822	2.1	ng/ul	163282	2	10.469	1531000	20.0
Tributyl phosphate	15.580	2.5	ng/ul	216410	3	14.333	1713540	20.0
n-Hexadecanoic ...	18.004	2.9	ng/ul	263240	4	17.092	1813530	20.0
Phenol, 4,4'-(1...	19.551	2.4	ng/ul	231023	5	21.292	1941540	20.0