

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM102824\
 Data File : BM048289.D
 Acq On : 29 Oct 2024 15:27
 Operator : RC/JU
 Sample : P4492-02
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4E61

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-BM102624.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BM048289.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.169	28	31	39	rVB	36653	49625	0.25%	0.013%
2	3.910	152	157	162	rBV	24932	37071	0.19%	0.010%
3	6.104	523	530	536	rBV4	29982	72937	0.37%	0.019%
4	6.922	656	669	684	rVB2	1094511	2723144	13.85%	0.719%
5	7.051	684	691	700	rBV	547924	855832	4.35%	0.226%
6	7.145	700	707	718	rVB	5026262	7629181	38.79%	2.014%
7	7.251	719	725	726	rBV	679581	814987	4.14%	0.215%
8	7.287	726	731	738	rVB	2834376	4624044	23.51%	1.221%
9	7.751	795	810	825	rBV	7205479	12558850	63.86%	3.316%
10	8.069	856	864	873	rVV	7190173	11680677	59.39%	3.084%
11	8.440	920	927	935	rBV	787297	1309938	6.66%	0.346%
12	8.534	935	943	957	rVB	5279579	8890749	45.21%	2.347%
13	8.792	979	987	994	rBV	4063424	6703570	34.09%	1.770%
14	8.875	994	1001	1004	rVV	610119	1045418	5.32%	0.276%
15	8.916	1004	1008	1017	rVB	1173499	1836303	9.34%	0.485%
16	9.292	1063	1072	1079	rBV2	38075	81604	0.41%	0.022%
17	9.451	1086	1099	1108	rBV	6235873	10738382	54.60%	2.835%
18	9.592	1116	1123	1125	rBV	521939	821994	4.18%	0.217%
19	9.622	1125	1128	1139	rVV	1131050	1893633	9.63%	0.500%
20	9.745	1142	1149	1158	rVB	261568	412034	2.10%	0.109%
21	9.928	1171	1180	1195	rBV	5833644	9509759	48.36%	2.511%
22	10.163	1209	1220	1235	rBV2	3868761	7644241	38.87%	2.018%
23	10.369	1251	1255	1263	rVB5	21538	38515	0.20%	0.010%
24	10.504	1269	1278	1285	rBV	856313	1499749	7.63%	0.396%
25	10.651	1295	1303	1318	rBV	357150	743597	3.78%	0.196%
26	10.839	1326	1335	1346	rVB	5676040	9363381	47.61%	2.472%
27	11.798	1491	1498	1517	rBV	1455813	2487911	12.65%	0.657%
28	12.063	1538	1543	1554	rVB2	61135	127045	0.65%	0.034%
29	12.175	1554	1562	1577	rBV	3351888	5538202	28.16%	1.462%
30	12.598	1628	1634	1639	rBV	78806	123765	0.63%	0.033%
31	12.792	1658	1667	1673	rBV	6753818	10459788	53.19%	2.762%
32	12.863	1673	1679	1696	rVB	1445518	2452885	12.47%	0.648%
33	13.186	1729	1734	1739	rVV	37980	49079	0.25%	0.013%
34	13.286	1747	1751	1757	rVV2	21397	35819	0.18%	0.009%
35	13.363	1757	1764	1768	rVV2	32126	67546	0.34%	0.018%
36	13.416	1768	1773	1780	rVB3	20338	40819	0.21%	0.011%

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Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : OFF Filtering: 5
 Sampling : 1 Min Area: 0.1 % of largest Peak
 Start Thrs: 0.2 Max Peaks: 100
 Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM102624.MA.M
 Title : SVOA CALIBRATION

37	13.774	1827	1834	1844	rBV	1270944	1889070	9.61%	0.499%
38	13.945	1855	1863	1872	rBV	6787442	10556370	53.68%	2.787%
39	14.051	1872	1881	1884	rVV	1572491	2637843	13.41%	0.696%
40	14.080	1884	1886	1893	rVB	1271956	1521404	7.74%	0.402%
41	14.363	1925	1934	1938	rBV	1136776	1756195	8.93%	0.464%
42	14.427	1938	1945	1951	rVB	7943507	12017832	61.11%	3.173%
43	14.486	1951	1955	1961	rVB	110965	147294	0.75%	0.039%
44	14.592	1964	1973	1983	rBV2	2258694	4066721	20.68%	1.074%
45	14.680	1983	1988	1991	rVV	95183	144162	0.73%	0.038%
46	14.739	1991	1998	2004	rVB	4815258	7071064	35.96%	1.867%
47	14.992	2036	2041	2052	rVB	105914	159466	0.81%	0.042%
48	15.221	2076	2080	2093	rVB5	38313	69728	0.35%	0.018%
49	15.357	2096	2103	2106	rBV	1854121	2839459	14.44%	0.750%
50	15.410	2106	2112	2119	rVV2	11454305	19666104	100.00%	5.192%
51	15.480	2119	2124	2136	rVV	500037	750545	3.82%	0.198%
52	15.657	2148	2154	2159	rVB2	177647	250750	1.28%	0.066%
53	15.721	2159	2165	2171	rBV	532553	745299	3.79%	0.197%
54	15.998	2208	2212	2217	rVB3	27434	35031	0.18%	0.009%
55	16.080	2221	2226	2234	rBV2	79543	150460	0.77%	0.040%
56	16.298	2253	2263	2270	rBV2	2156494	3157755	16.06%	0.834%
57	16.415	2275	2283	2290	rBV	9697138	15075063	76.66%	3.980%
58	16.762	2330	2342	2351	rBV	7842909	11602567	59.00%	3.063%
59	16.904	2362	2366	2375	rVB9	51125	124063	0.63%	0.033%
60	17.104	2393	2400	2403	rBV	1249136	1866229	9.49%	0.493%
61	17.145	2403	2407	2412	rVV	3621154	4969005	25.27%	1.312%
62	17.198	2412	2416	2419	rVV	1895282	2831546	14.40%	0.748%
63	17.233	2419	2422	2436	rVB	3377662	4578832	23.28%	1.209%
64	17.392	2441	2449	2454	rVB4	28567	64177	0.33%	0.017%
65	17.515	2465	2470	2475	rVB3	36063	54662	0.28%	0.014%
66	17.651	2489	2493	2499	rVV8	15950	40782	0.21%	0.011%
67	17.715	2499	2504	2510	rVV4	70020	134375	0.68%	0.035%
68	17.774	2510	2514	2522	rVB	173852	222302	1.13%	0.059%
69	17.998	2548	2552	2558	rBV	363389	506928	2.58%	0.134%
70	18.074	2558	2565	2570	rVV	10773002	14234609	72.38%	3.758%
71	18.168	2578	2581	2586	rVB5	25134	35414	0.18%	0.009%
72	18.298	2597	2603	2607	rBV3	54653	97623	0.50%	0.026%
73	18.368	2611	2615	2621	rBV6	48679	97095	0.49%	0.026%
74	18.457	2626	2630	2637	rVV4	29404	59807	0.30%	0.016%
75	18.521	2637	2641	2648	rVV	317586	435957	2.22%	0.115%
76	18.715	2670	2674	2678	rVB6	31641	41727	0.21%	0.011%

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Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

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77	18.845	2691	2696	2703	rVB	84937	149382	0.76%	0.039%
78	18.951	2711	2714	2717	rVB3	37437	40306	0.20%	0.011%
79	19.056	2729	2732	2734	rBV2	27430	36476	0.19%	0.010%
80	19.086	2734	2737	2741	rVV2	59319	82433	0.42%	0.022%
81	19.162	2743	2750	2762	rVB	11324106	16275036	82.76%	4.297%
82	19.274	2766	2769	2774	rBV2	58427	87066	0.44%	0.023%
83	19.404	2787	2791	2796	rBV	106328	135810	0.69%	0.036%
84	19.492	2800	2806	2808	rBV	2462965	3476462	17.68%	0.918%
85	19.521	2808	2811	2820	rVB	5618910	7612083	38.71%	2.010%
86	19.645	2828	2832	2836	rBV5	47780	62302	0.32%	0.016%
87	20.180	2920	2923	2927	rBV2	176868	195997	1.00%	0.052%
88	20.421	2959	2964	2970	rVB	3116173	3654783	18.58%	0.965%
89	21.009	3060	3064	3074	rVB3	464407	756699	3.85%	0.200%
90	21.156	3085	3089	3095	rBV2	172684	231393	1.18%	0.061%
91	21.233	3096	3102	3108	rVV	12310131	16797549	85.41%	4.435%
92	21.309	3109	3115	3121	rVV	8593691	15536740	79.00%	4.102%
93	21.380	3121	3127	3138	rVB	11943992	19298224	98.13%	5.095%
94	22.256	3273	3276	3288	rVB5	139630	326366	1.66%	0.086%
95	23.356	3452	3463	3470	rBV	7913878	18648179	94.82%	4.924%
96	24.068	3576	3584	3589	rBV	1071398	2738115	13.92%	0.723%
97	24.133	3589	3595	3607	rVB	1790298	4287242	21.80%	1.132%
98	24.268	3611	3618	3636	rVB	726771	1840408	9.36%	0.486%
99	27.644	4173	4192	4211	rVB2	2505193	11733297	59.66%	3.098%
100	28.615	4347	4357	4375	rVB2	724839	3079467	15.66%	0.813%

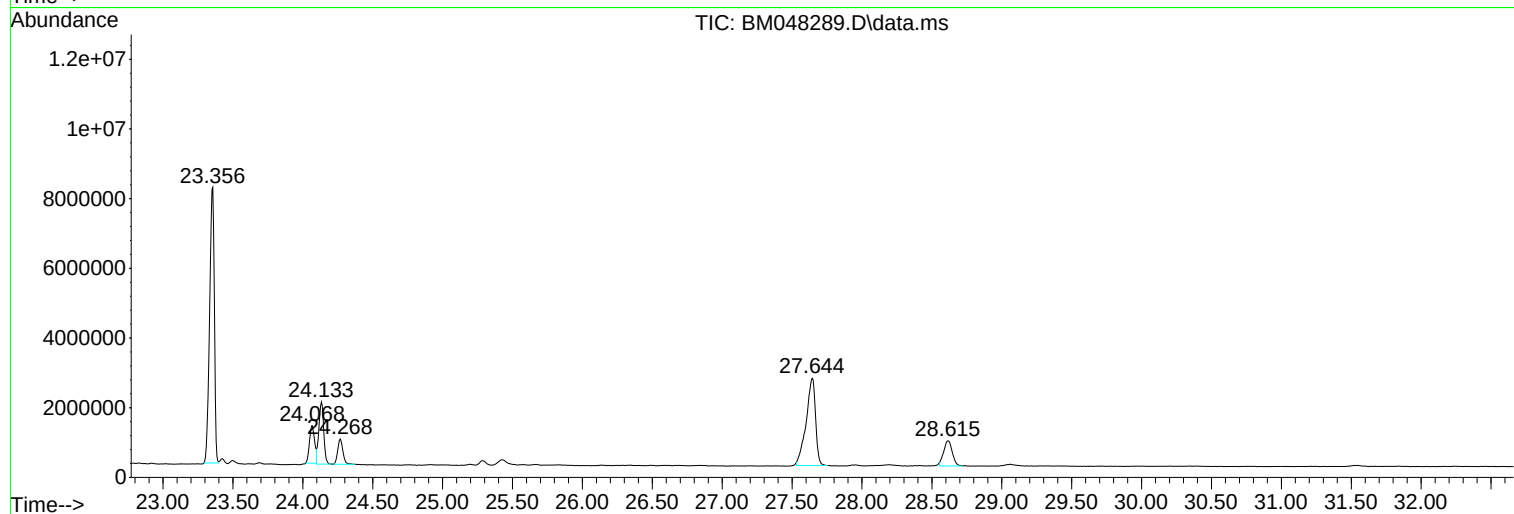
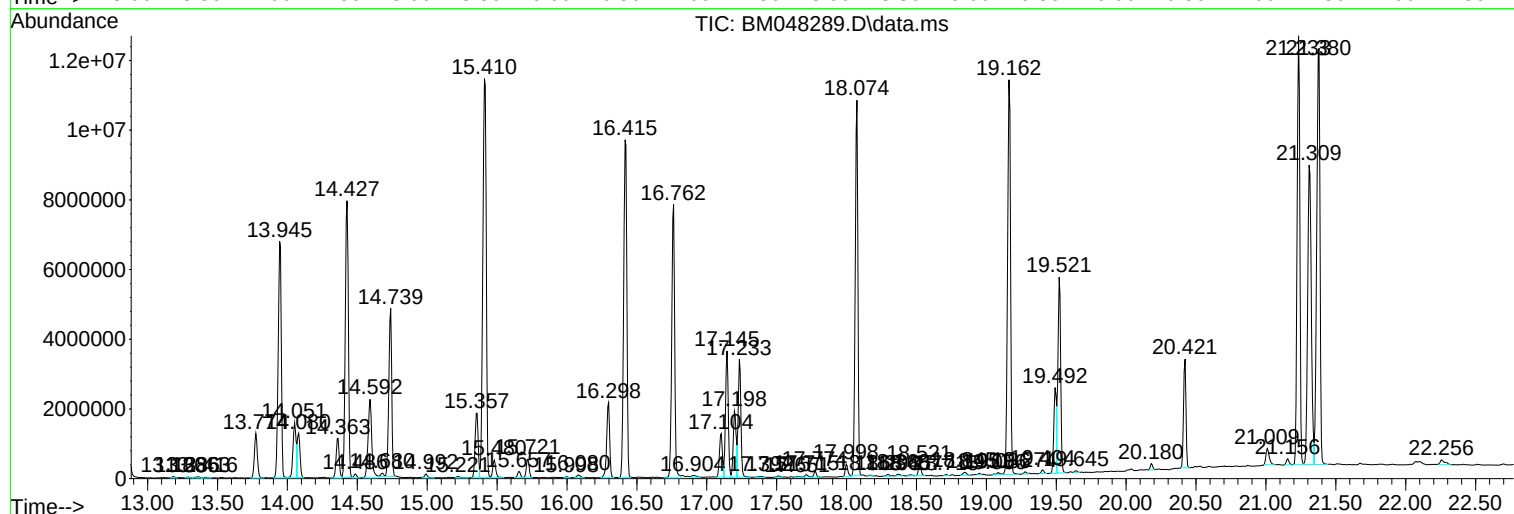
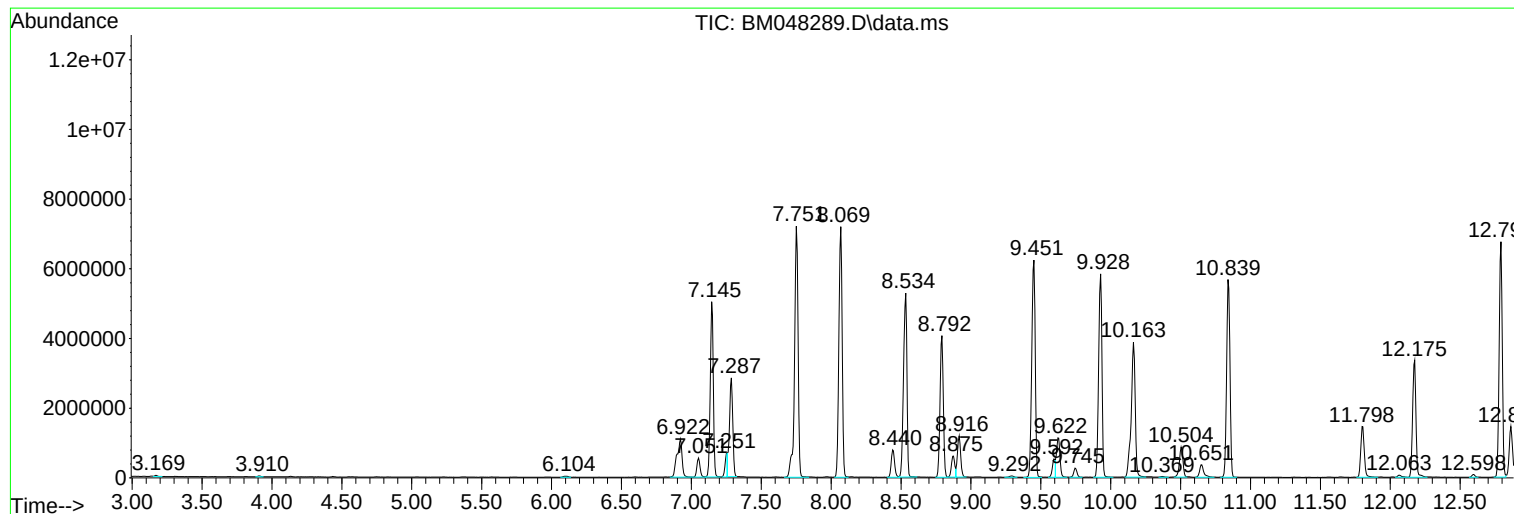
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Instrument :
BNA_M
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM102624.MA.M
Quant Title : SVOA CALIBRATION

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



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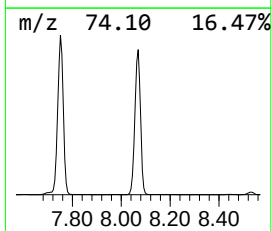
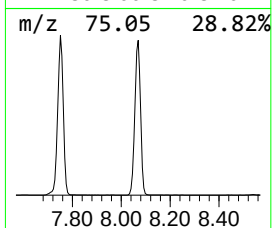
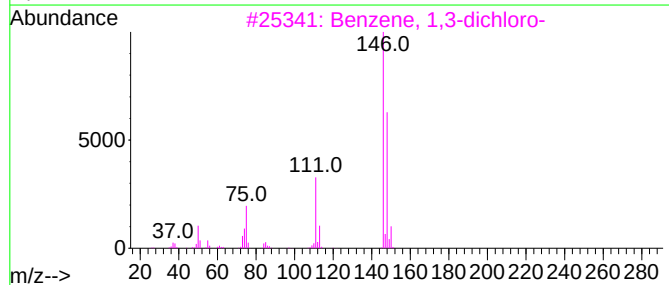
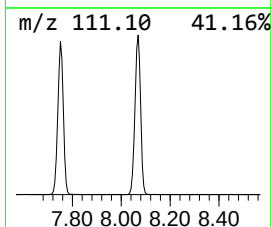
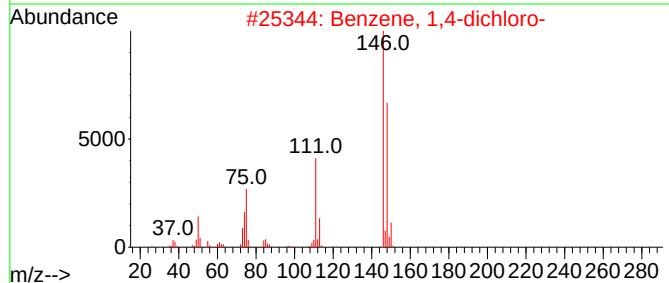
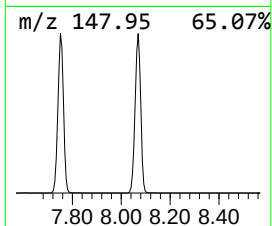
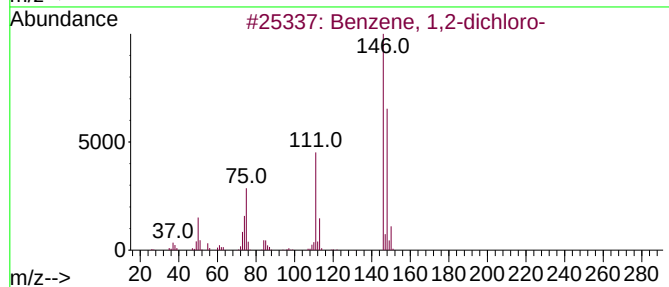
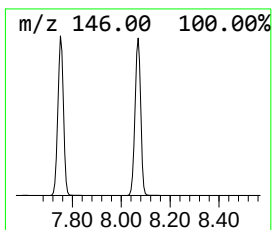
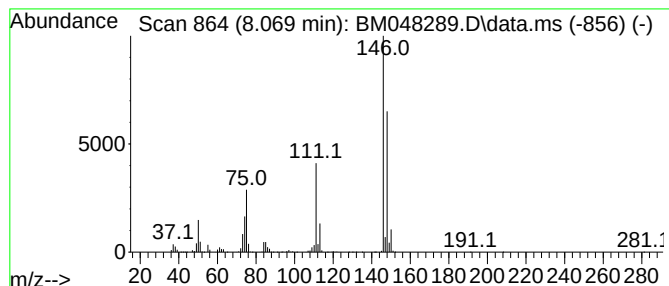
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM102624.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzene, 1,2-dichloro- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.069	18.60 ng/ul	11680700	1,4-Dichlorobenzene-d4	7.716

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	98
2			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	98
3			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	97
4			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	97
5			Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	97



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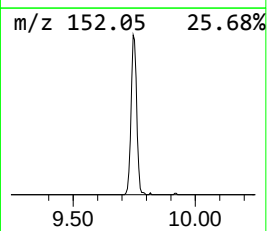
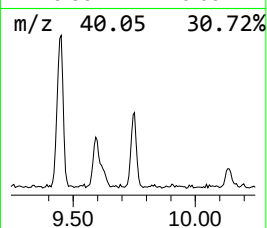
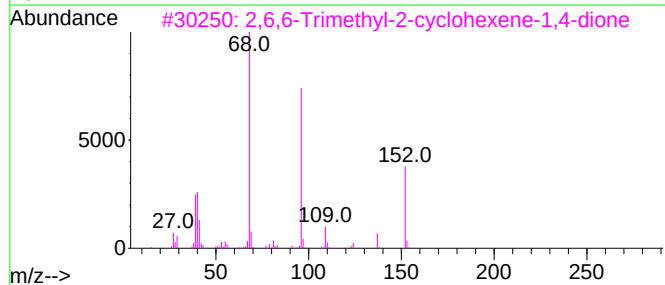
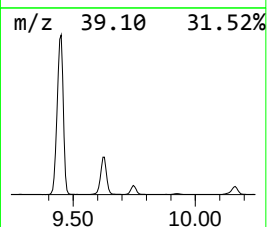
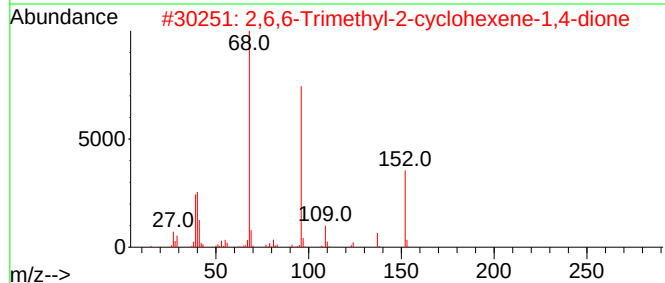
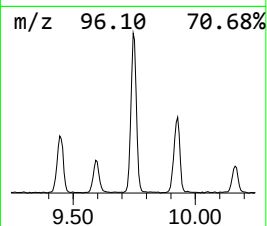
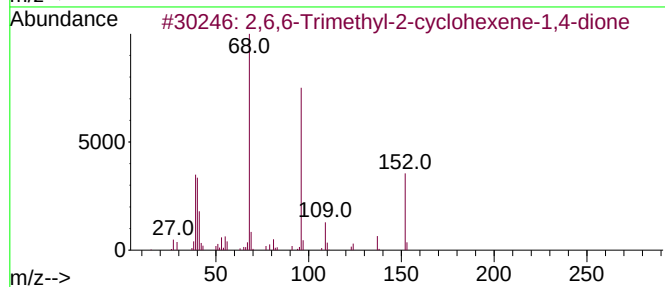
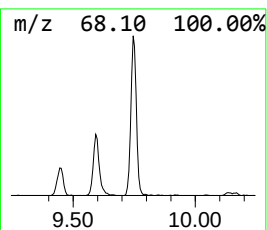
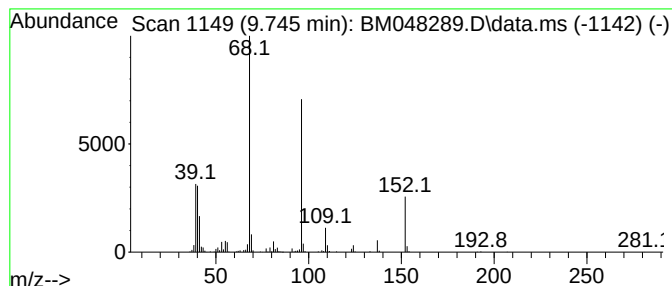
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM102624.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 2,6,6-Trimethyl-2-cyclohexe... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.745	5.49 ng/ul	412034	Naphthalene-d8	10.504

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,6,6-Trimethyl-2-cyclohexene-1,...	152	C9H12O2	001125-21-9	96		
2	2,6,6-Trimethyl-2-cyclohexene-1,...	152	C9H12O2	001125-21-9	95		
3	2,6,6-Trimethyl-2-cyclohexene-1,...	152	C9H12O2	001125-21-9	94		
4	2,6,6-Trimethyl-2-cyclohexene-1,...	152	C9H12O2	001125-21-9	91		
5	2-Cyclohexen-1-one	96	C6H8O	000930-68-7	53		



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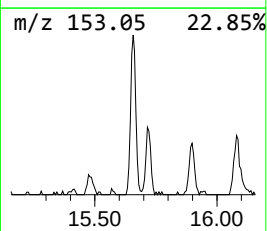
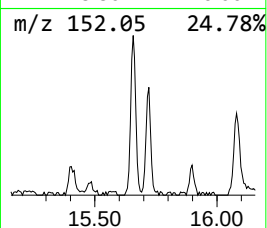
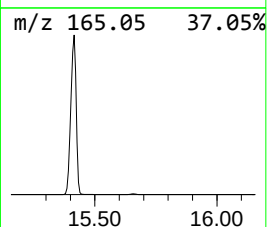
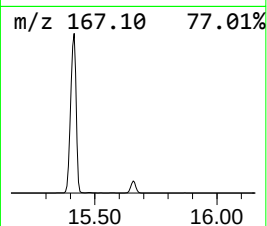
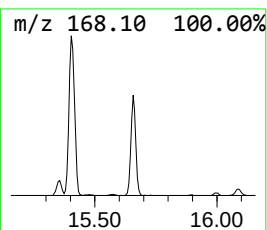
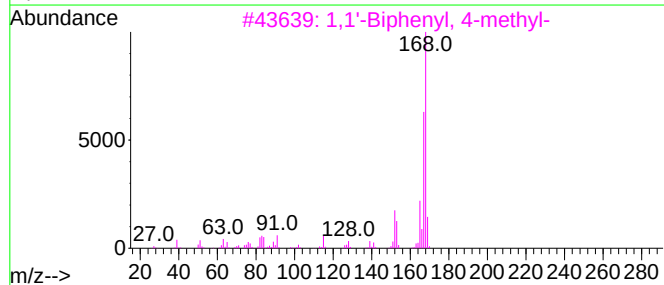
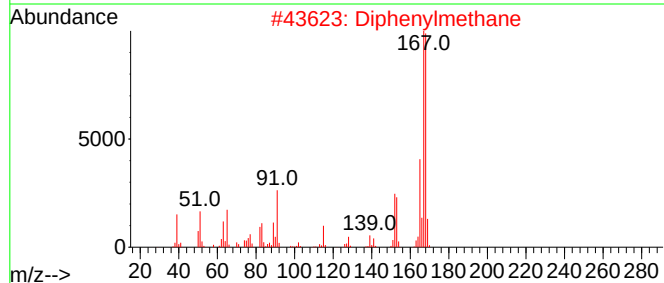
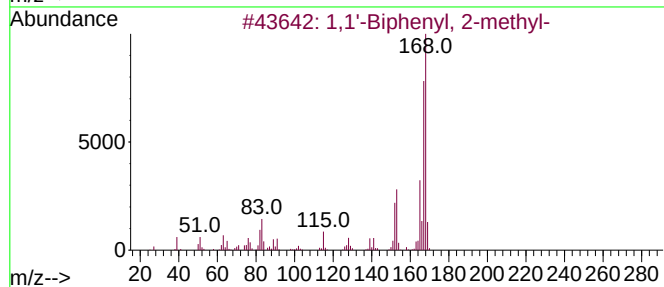
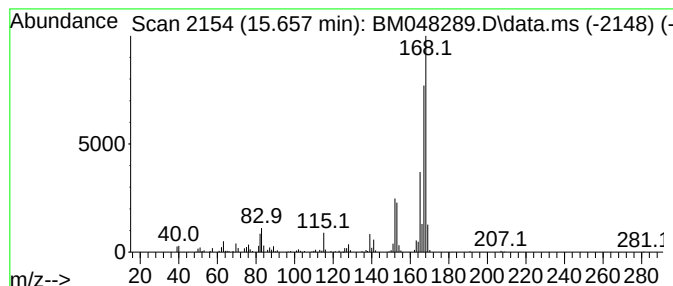
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 1,1'-Biphenyl, 2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.657	2.86 ng/ul	250750	Acenaphthene-d10	14.363

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1,1'-Biphenyl, 2-methyl-	168	C13H12	000643-58-3	94
2		Diphenylmethane	168	C13H12	000101-81-5	93
3		1,1'-Biphenyl, 4-methyl-	168	C13H12	000644-08-6	93
4		Diphenylmethane	168	C13H12	000101-81-5	93
5		1,1'-Biphenyl, 2-methyl-	168	C13H12	000643-58-3	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM102824\
Data File : BM048289.D
Acq On : 29 Oct 2024 15:27
Operator : RC/JU
Sample : P4492-02
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
A4E61

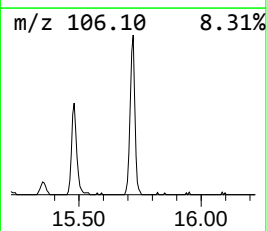
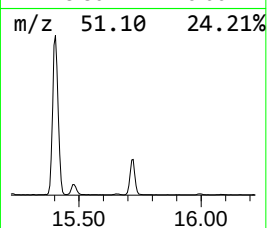
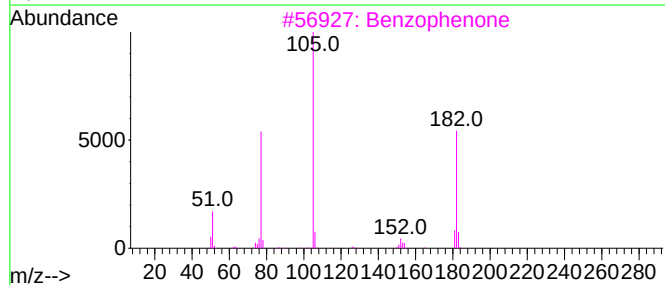
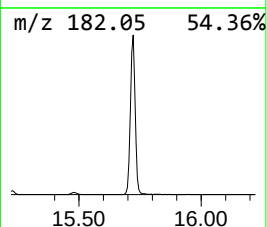
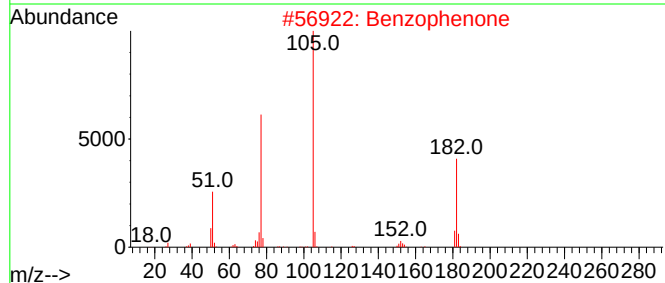
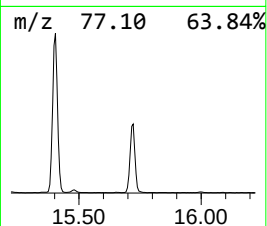
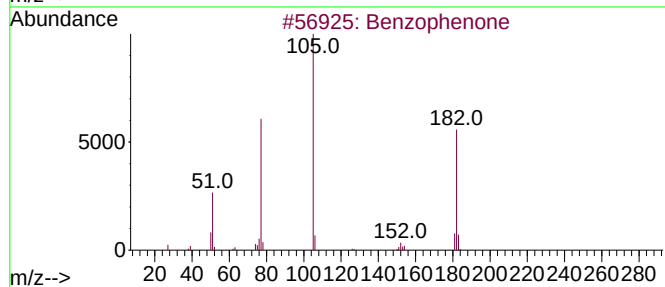
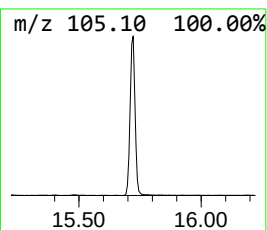
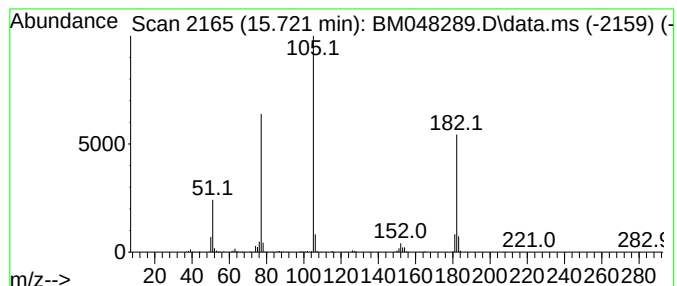
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Benzophenone Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.721	8.49 ng/ul	745299	Acenaphthene-d10	14.363

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	96
2			Benzophenone	182	C13H10O	000119-61-9	96
3			Benzophenone	182	C13H10O	000119-61-9	95
4			Benzophenone	182	C13H10O	000119-61-9	93
5			Benzenecarbothioic acid, S-methy...	152	C8H8OS	005925-68-8	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM102824\
Data File : BM048289.D
Acq On : 29 Oct 2024 15:27
Operator : RC/JU
Sample : P4492-02
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
A4E61

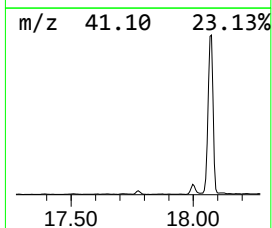
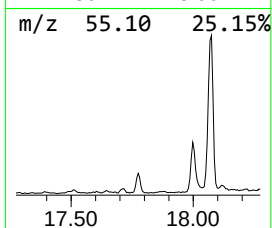
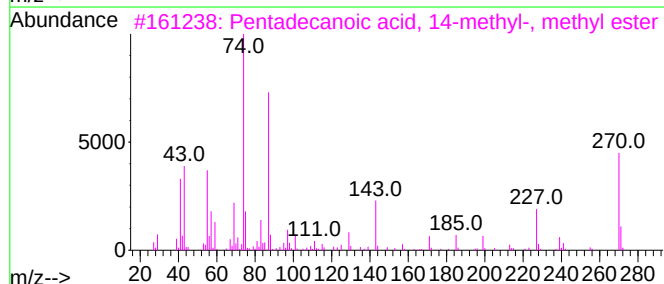
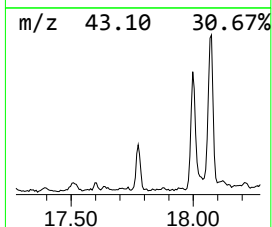
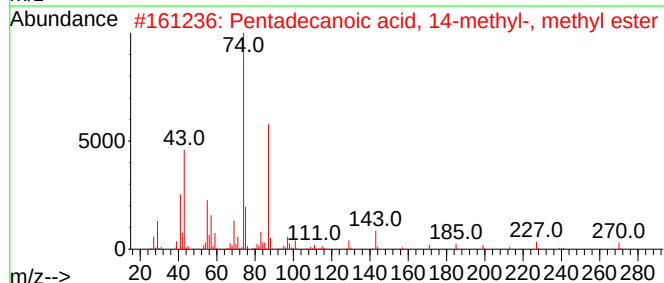
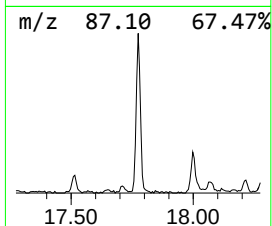
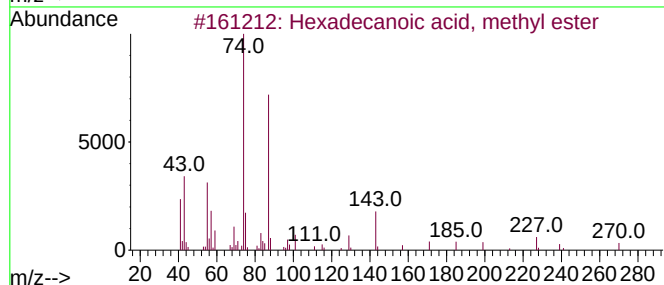
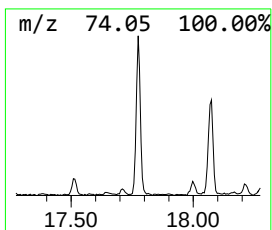
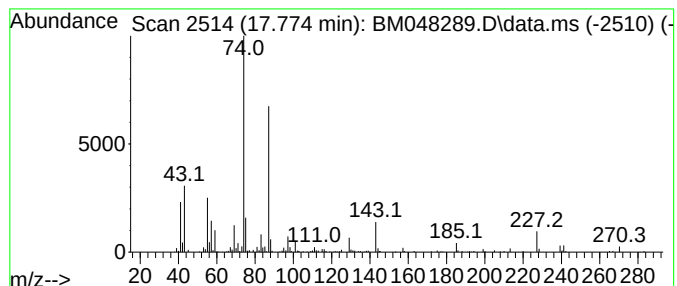
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Hexadecanoic acid, methyl e... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.774	2.38 ng/ul	222302	Phenanthrene-d10	17.104

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	95
2			Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	93
3			Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	91
4			Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	91
5			Pentadecanoic acid, methyl ester	256	C16H32O2	007132-64-1	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM102824\
Data File : BM048289.D
Acq On : 29 Oct 2024 15:27
Operator : RC/JU
Sample : P4492-02
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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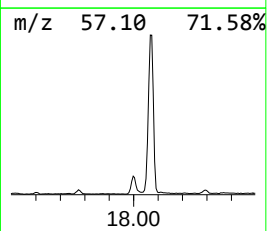
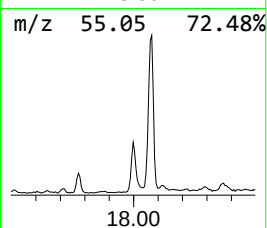
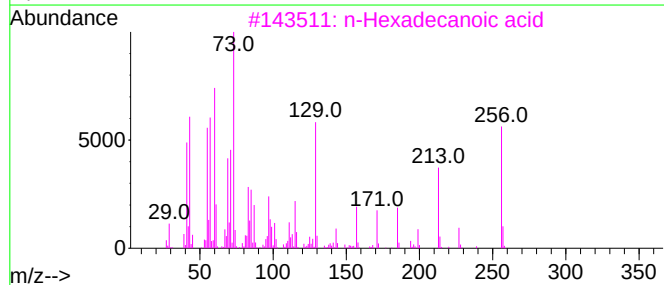
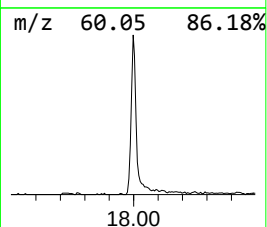
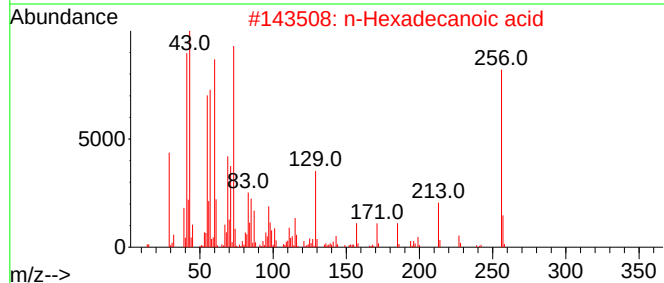
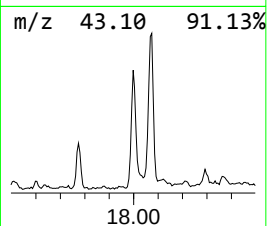
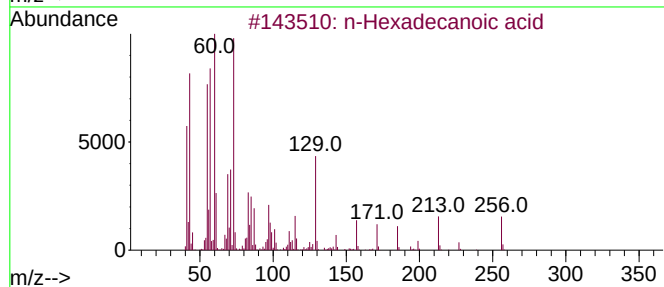
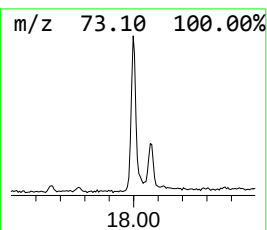
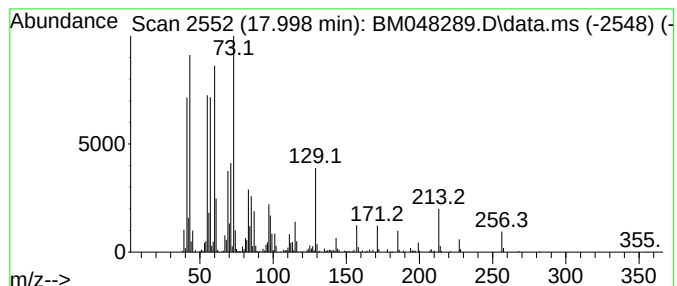
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.998	5.43 ng/ul	506928	Phenanthrene-d10	17.104

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	99
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
5			Tridecanoic acid	214	C13H26O2	000638-53-9	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM102824\
Data File : BM048289.D
Acq On : 29 Oct 2024 15:27
Operator : RC/JU
Sample : P4492-02
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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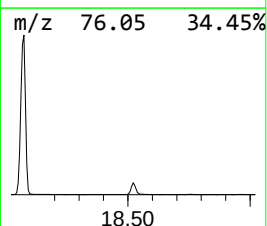
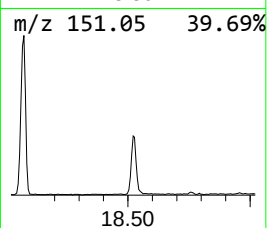
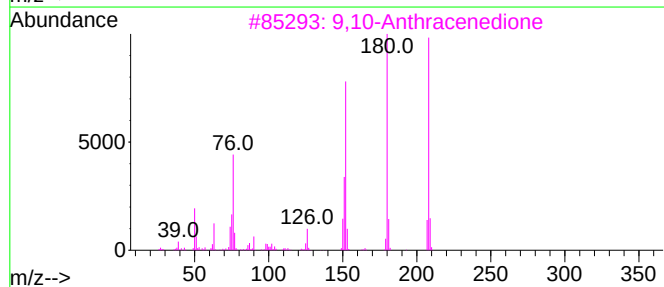
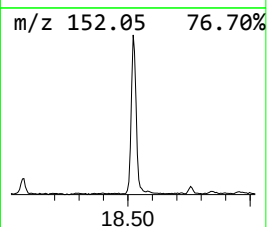
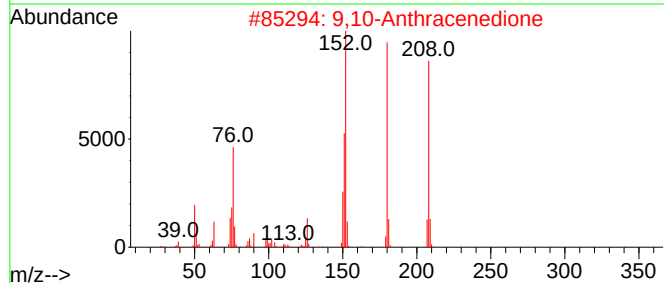
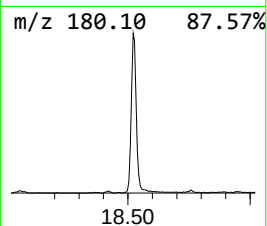
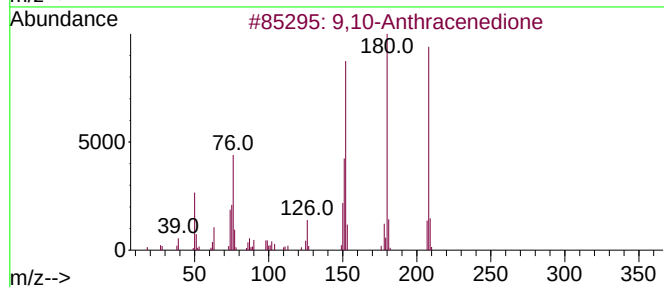
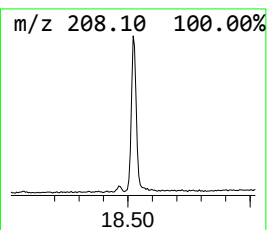
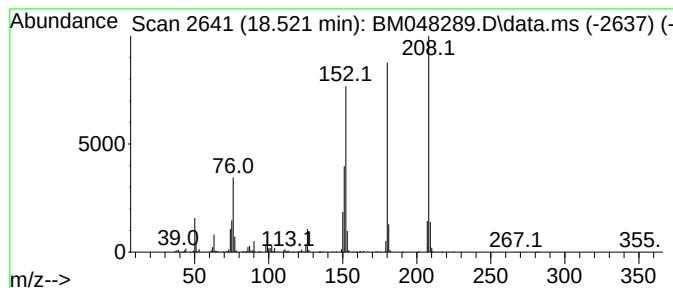
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 9,10-Anthracenedione Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.521	4.67 ng/ul	435957	Phenanthrene-d10	17.104

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,10-Anthracenedione			208	C14H8O2	000084-65-1	99
2	9,10-Anthracenedione			208	C14H8O2	000084-65-1	98
3	9,10-Anthracenedione			208	C14H8O2	000084-65-1	97
4	9,10-Anthracenedione			208	C14H8O2	000084-65-1	96
5	9H-Fluoren-9-one			180	C13H8O	000486-25-9	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM102824\
Data File : BM048289.D
Acq On : 29 Oct 2024 15:27
Operator : RC/JU
Sample : P4492-02
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM102624.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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2,6,6-Trimethyl...	9.745	5.5	ng/ul	412034	2	10.504	1499750	20.0
1,1'-Biphenyl, ...	15.657	2.9	ng/ul	250750	3	14.363	1756200	20.0
Benzophenone	15.721	8.5	ng/ul	745299	3	14.363	1756200	20.0
Hexadecanoic ac...	17.774	2.4	ng/ul	222302	4	17.104	1866230	20.0
n-Hexadecanoic ...	17.998	5.4	ng/ul	506928	4	17.104	1866230	20.0
9,10-Anthracene...	18.521	4.7	ng/ul	435957	4	17.104	1866230	20.0