

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061224\
 Data File : BN031913.D
 Acq On : 11 Jun 2024 16:28
 Operator : MA/JU
 Sample : P2642-18
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A4DG2

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_N\Methods\SFAM-EPA-BN052924.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BN031913.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.175	29	32	40	rVB	54343	79666	0.19%	0.011%
2	3.604	103	105	122	rBV3	53256	132426	0.32%	0.018%
3	6.040	508	519	522	rBV	49793	76170	0.19%	0.010%
4	6.869	647	660	677	rVB2	6179009	12093057	29.52%	1.618%
5	7.004	677	683	693	rBV	1011388	1579658	3.86%	0.211%
6	7.098	693	699	709	rVB	1861650	2755133	6.73%	0.369%
7	7.198	709	716	717	rBV	1207754	1782337	4.35%	0.239%
8	7.228	717	721	728	rVB	4128571	6628444	16.18%	0.887%
9	7.545	768	775	784	rBV	2348712	3853051	9.41%	0.516%
10	7.698	795	801	813	rVB	8519372	14282299	34.87%	1.911%
11	8.010	845	854	867	rVB	6199876	10322405	25.20%	1.381%
12	8.381	909	917	925	rBV	1543126	2544651	6.21%	0.341%
13	8.475	925	933	942	rBV	5179999	8871334	21.66%	1.187%
14	8.728	964	976	984	rBV	5220645	8793732	21.47%	1.177%
15	8.816	984	991	995	rBV	1151811	2055098	5.02%	0.275%
16	8.863	995	999	1006	rVB	2120437	3301502	8.06%	0.442%
17	9.222	1053	1060	1066	rVB3	54537	96351	0.24%	0.013%
18	9.534	1106	1113	1114	rBV	965625	1273272	3.11%	0.170%
19	9.569	1114	1119	1133	rVB	2342958	4220919	10.30%	0.565%
20	9.869	1162	1170	1180	rVB	3820637	6112832	14.92%	0.818%
21	10.098	1197	1209	1223	rVB2	3551502	8414178	20.54%	1.126%
22	10.445	1260	1268	1271	rVV	1498831	2595050	6.34%	0.347%
23	10.498	1271	1277	1285	rVV	11989449	21695689	52.97%	2.904%
24	10.586	1285	1292	1302	rVV	461066	917655	2.24%	0.123%
25	11.022	1358	1366	1370	rBV	142853	251428	0.61%	0.034%
26	11.075	1370	1375	1389	rVV	153868	278347	0.68%	0.037%
27	11.192	1389	1395	1416	rVB	142126	328593	0.80%	0.044%
28	11.739	1474	1488	1503	rBV	6875835	11742534	28.67%	1.572%
29	12.545	1617	1625	1636	rVB	151476	229952	0.56%	0.031%
30	12.727	1649	1656	1662	rVV	4966918	8292964	20.25%	1.110%
31	12.804	1662	1669	1683	rVV	10096105	16845402	41.12%	2.254%
32	12.939	1683	1692	1695	rVV4	254958	580344	1.42%	0.078%
33	12.969	1695	1697	1706	rVB	245801	338621	0.83%	0.045%
34	13.175	1726	1732	1739	rVV	7026468	11385153	27.79%	1.524%
35	13.233	1739	1742	1752	rVB2	55644	91898	0.22%	0.012%
36	13.357	1759	1763	1772	rVB	99216	167229	0.41%	0.022%

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Integration Parameters: LSCINT.P

Integrator: RTE

Soothing : 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_N\Methods\SFAM-EPA-BN052924.MA.M
Title : SVOA CALIBRATION

37	13.727	1819	1826	1835	rBV	3017707	4678909	11.42%	0.626%
38	13.910	1847	1857	1862	rBV	12305768	19920363	48.63%	2.666%
39	14.033	1862	1878	1885	rVV2	15291246	30990155	75.66%	4.147%
40	14.310	1917	1925	1930	rBV	2362979	3602593	8.80%	0.482%
41	14.374	1930	1936	1942	rVV	13907411	22562769	55.08%	3.020%
42	14.433	1942	1946	1954	rVB	267842	379205	0.93%	0.051%
43	14.563	1954	1968	1974	rBV2	11636820	26402877	64.46%	3.534%
44	14.622	1974	1978	1983	rVV	288971	564272	1.38%	0.076%
45	14.698	1983	1991	1996	rVV	7678896	12014767	29.33%	1.608%
46	14.733	1996	1997	2008	rVV2	90661	171397	0.42%	0.023%
47	14.939	2027	2032	2044	rVB	208459	327006	0.80%	0.044%
48	15.151	2056	2068	2073	rBV	16280832	26549734	64.82%	3.553%
49	15.304	2087	2094	2098	rBV	4643528	6835207	16.69%	0.915%
50	15.363	2098	2104	2111	rVV	16795545	27063633	66.07%	3.622%
51	15.439	2111	2117	2128	rVB	1410383	2035217	4.97%	0.272%
52	15.610	2142	2146	2151	rBV	113641	149468	0.36%	0.020%
53	15.857	2182	2188	2189	rBV	49710	77315	0.19%	0.010%
54	15.898	2191	2195	2200	rVV	142037	258972	0.63%	0.035%
55	15.939	2200	2202	2212	rVB3	54007	100301	0.24%	0.013%
56	16.039	2212	2219	2226	rBV3	156480	310087	0.76%	0.041%
57	16.221	2243	2250	2261	rVB	235739	321333	0.78%	0.043%
58	16.368	2266	2275	2281	rBV	16129819	29130577	71.12%	3.899%
59	16.716	2326	2334	2343	rBV	13886963	22352547	54.57%	2.991%
60	16.863	2354	2359	2372	rVB6	162780	396435	0.97%	0.053%
61	17.057	2384	2392	2395	rVV	2921775	4351797	10.62%	0.582%
62	17.104	2395	2400	2404	rVV	9436407	14069891	34.35%	1.883%
63	17.157	2404	2409	2412	rVV2	4921342	8057519	19.67%	1.078%
64	17.198	2412	2416	2428	rVB	11598442	16753592	40.90%	2.242%
65	17.333	2433	2439	2451	rVB3	52758	149300	0.36%	0.020%
66	17.468	2456	2462	2467	rVB	154010	229800	0.56%	0.031%
67	17.604	2480	2485	2488	rVV3	66223	108900	0.27%	0.015%
68	17.668	2492	2496	2502	rVV2	109310	197991	0.48%	0.026%
69	17.733	2502	2507	2518	rVB	787967	1018428	2.49%	0.136%
70	17.910	2533	2537	2541	rBV5	62443	90767	0.22%	0.012%
71	17.963	2541	2546	2552	rBV	985497	1478258	3.61%	0.198%
72	18.039	2552	2559	2564	rVB	17820438	28272422	69.02%	3.784%
73	18.121	2569	2573	2578	rVB4	108116	173874	0.42%	0.023%
74	18.251	2590	2595	2598	rBV3	80031	138410	0.34%	0.019%
75	18.327	2604	2608	2613	rBV3	217935	380412	0.93%	0.051%
76	18.421	2620	2624	2629	rBV	129944	175293	0.43%	0.023%

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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_N\Methods\SFAM-EPA-BN052924.MA.M
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77	18.480	2629	2634	2644	rVB	1596867	2235446	5.46%	0.299%
78	18.839	2690	2695	2699	rBV4	83734	177236	0.43%	0.024%
79	18.880	2700	2702	2704	rVV2	122245	142919	0.35%	0.019%
80	18.915	2704	2708	2711	rVV	735236	883093	2.16%	0.118%
81	18.945	2711	2713	2720	rVB5	105135	147082	0.36%	0.020%
82	19.015	2722	2725	2727	rBV2	94169	113374	0.28%	0.015%
83	19.051	2727	2731	2736	rVV2	323845	391916	0.96%	0.052%
84	19.133	2736	2745	2750	rVV	19462800	34838206	85.05%	4.662%
85	19.245	2760	2764	2779	rBV	499799	934084	2.28%	0.125%
86	19.368	2780	2785	2793	rVV	326128	522050	1.27%	0.070%
87	19.462	2794	2801	2802	rVV	6471164	9180873	22.41%	1.229%
88	19.492	2802	2806	2814	rVB	13969075	20796918	50.77%	2.783%
89	20.151	2915	2918	2922	rBV	448196	503413	1.23%	0.067%
90	20.992	3056	3061	3071	rVB2	572393	1063091	2.60%	0.142%
91	21.133	3081	3085	3090	rBV	827877	968164	2.36%	0.130%
92	21.215	3092	3099	3104	rVV	21604732	36438985	88.96%	4.877%
93	21.286	3104	3111	3117	rVV2	15979571	33645227	82.14%	4.503%
94	21.356	3117	3123	3133	rVB	19162162	34919043	85.25%	4.673%
95	22.045	3236	3240	3243	rBV3	284049	433862	1.06%	0.058%
96	22.233	3267	3272	3280	rVB	610722	1042374	2.54%	0.140%
97	23.386	3455	3468	3477	rVB	14615262	40961675	100.00%	5.482%
98	24.027	3568	3577	3588	rVB	3488988	8521467	20.80%	1.140%
99	24.227	3603	3611	3622	rVB2	1880454	4769421	11.64%	0.638%
100	27.603	4168	4185	4202	rVB2	6108780	25732915	62.82%	3.444%

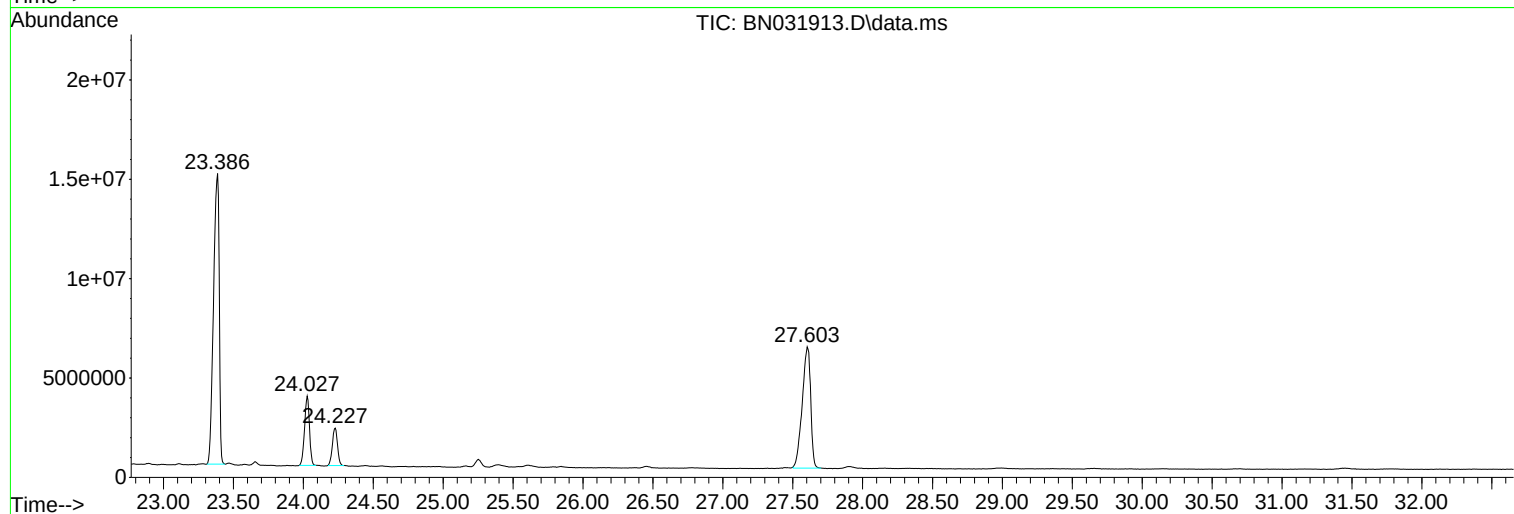
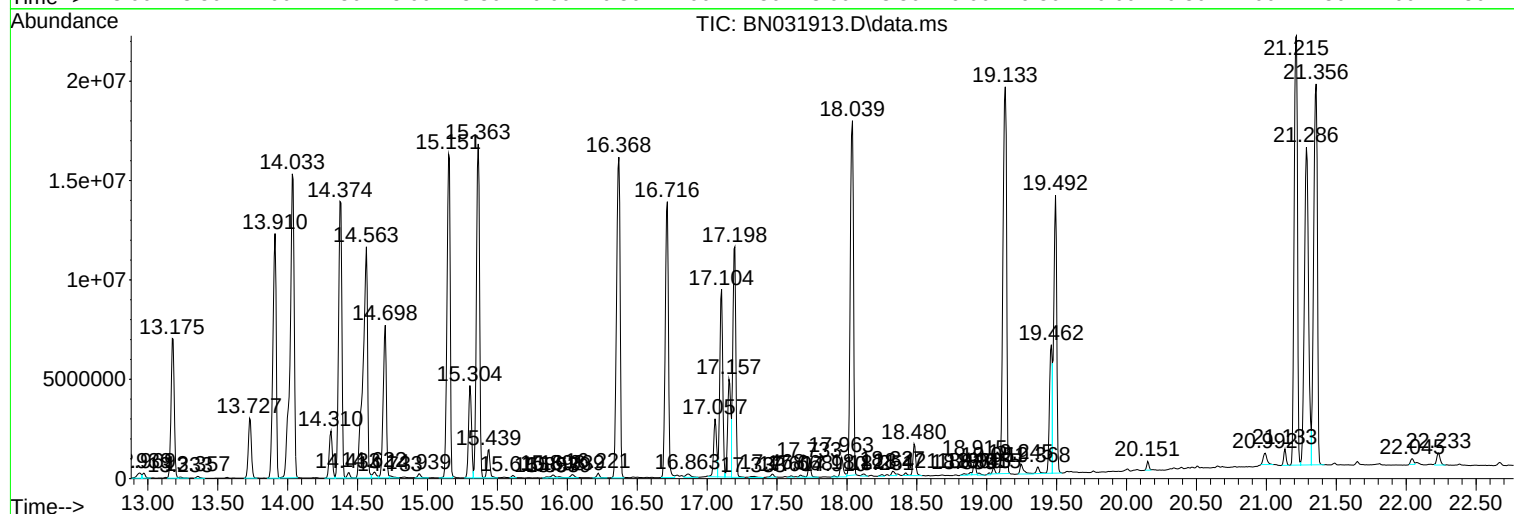
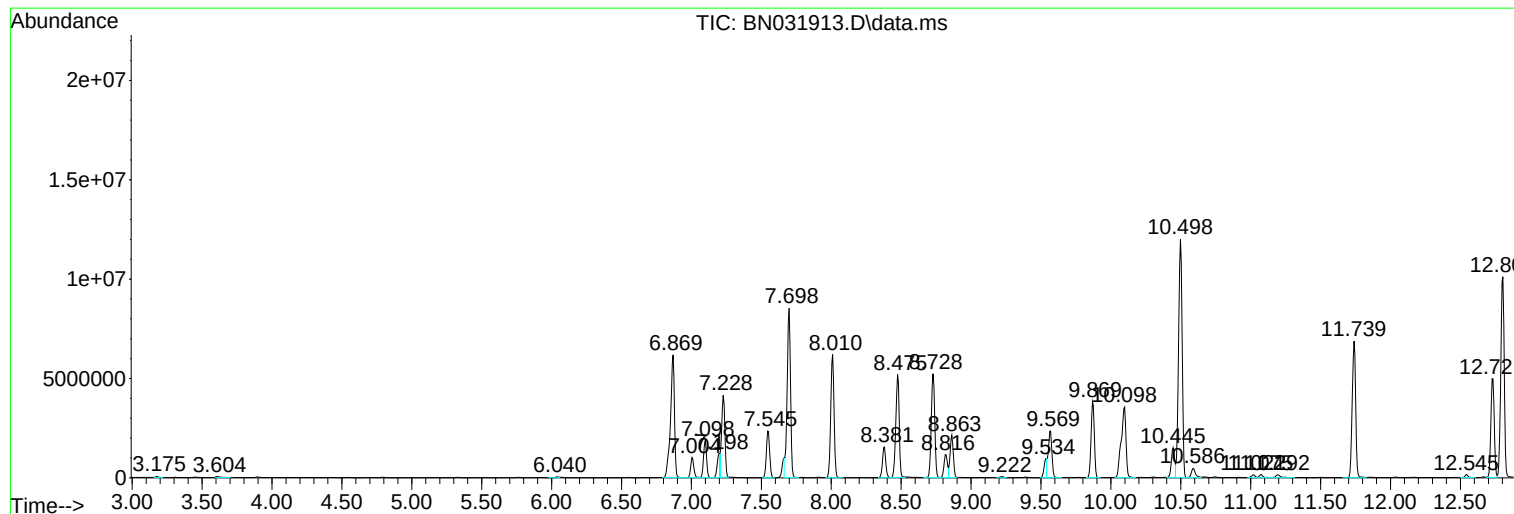
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Instrument :
BNA_N
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

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



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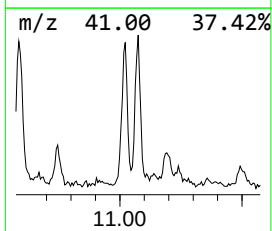
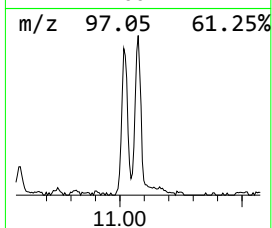
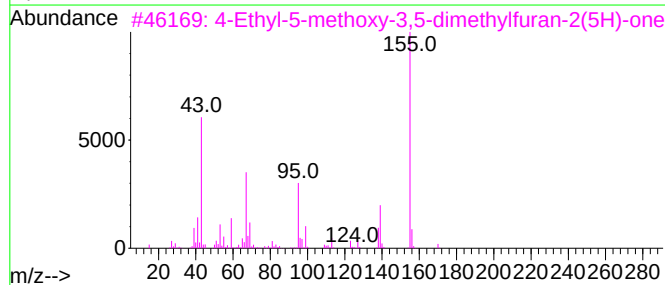
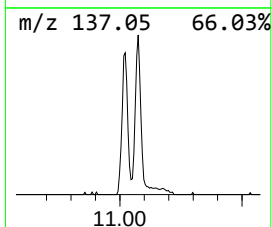
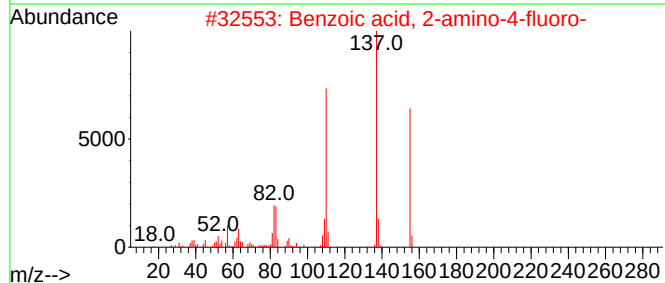
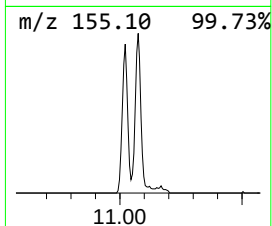
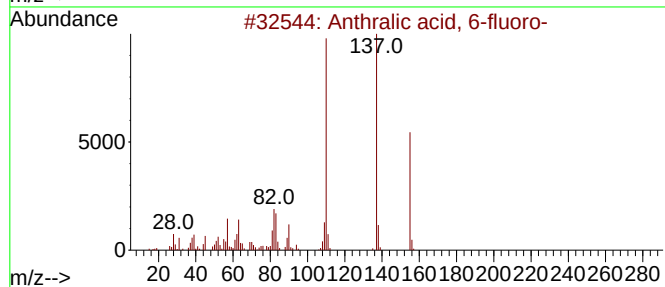
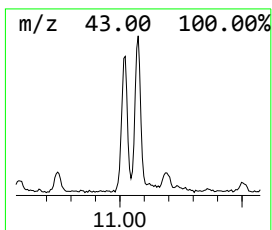
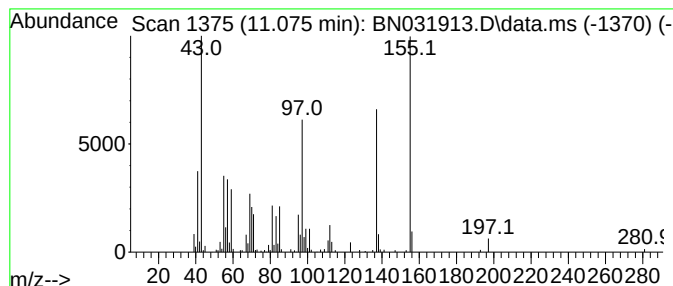
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 unknown-01 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.075	2.15 ng/ul	278347	Naphthalene-d8	10.445

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Anthralic acid, 6-fluoro-	155	C7H6FN02	000434-76-4	38
2			Benzoic acid, 2-amino-4-fluoro-	155	C7H6FN02	000446-32-2	38
3			4-Ethyl-5-methoxy-3,5-dimethylfu...	170	C9H14O3	1000495-96-7	35
4			2-Methoxyethyl pinacolyl methylp...	238	C10H23O4P	177600-27-0	35
5			2-Amino-4,6-dimethoxypyrimidine	155	C6H9N3O2	036315-01-2	27



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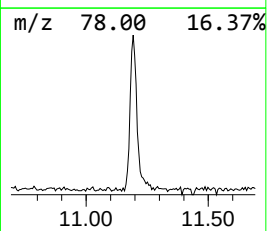
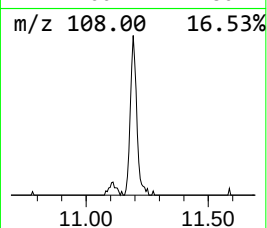
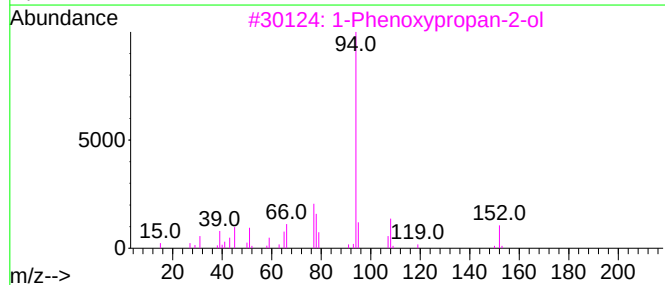
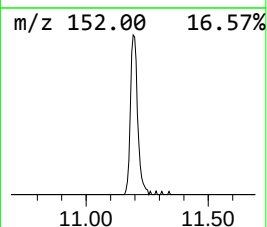
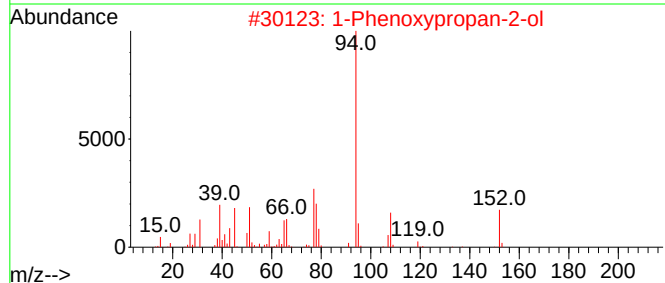
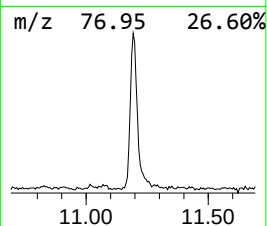
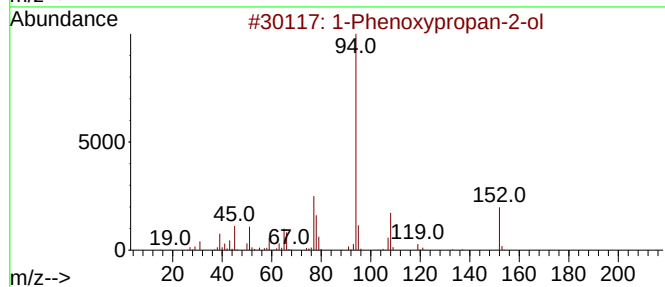
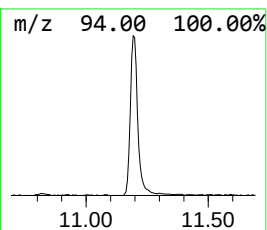
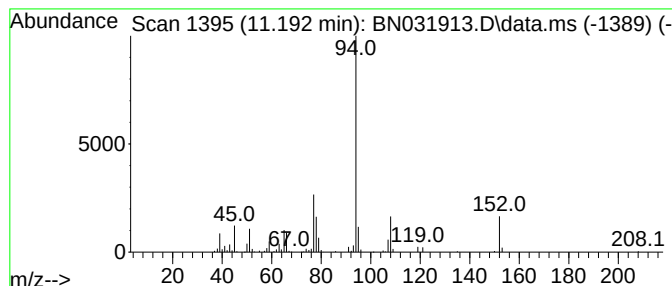
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 4 1-Phenoxypropan-2-ol Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.192	2.53 ng/ul	328593	Naphthalene-d8	10.445

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	96
2			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	95
3			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	93
4			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	87
5			Ethanol, 2-phenoxy-	138	C8H10O2	000122-99-6	64



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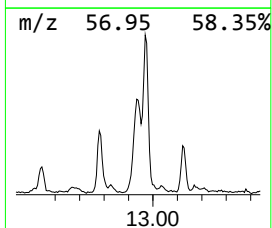
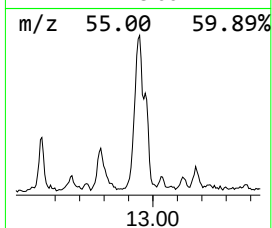
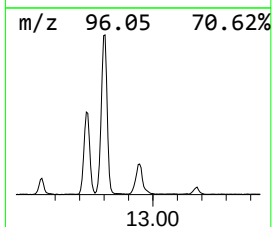
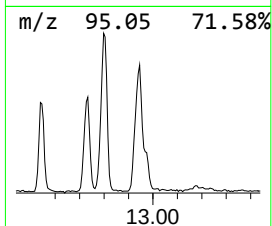
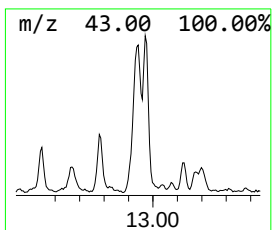
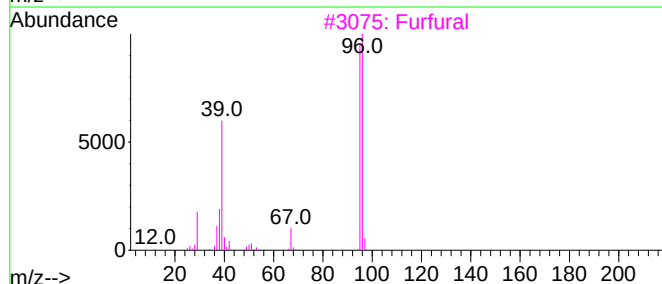
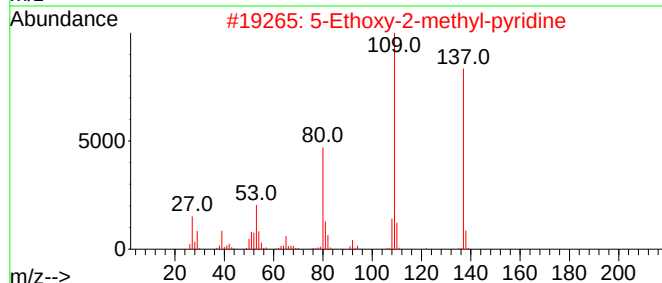
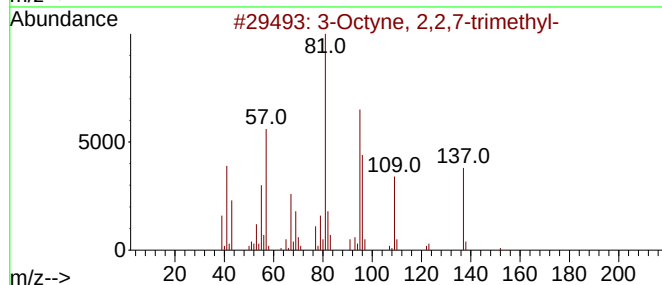
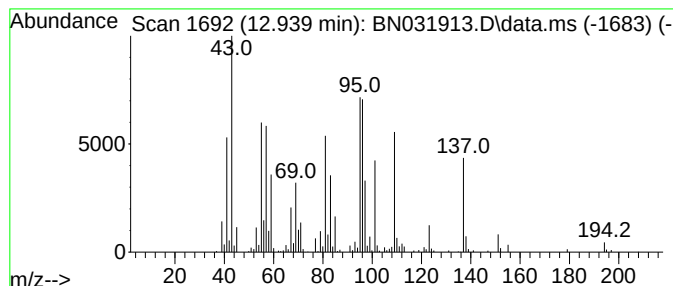
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 unknown-02 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.939	3.22 ng/ul	580344	Acenaphthene-d10	14.310

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Octyne, 2,2,7-trimethyl-	152	C11H20	055402-13-6	49
2			5-Ethoxy-2-methyl-pyridine	137	C8H11NO	055270-48-9	18
3			Furfural	96	C5H4O2	000098-01-1	18
4			Furfural	96	C5H4O2	000098-01-1	18
5			Furfural	96	C5H4O2	000098-01-1	18



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061224\
Data File : BN031913.D
Acq On : 11 Jun 2024 16:28
Operator : MA/JU
Sample : P2642-18
Misc :
ALS Vial : 4 Sample Multiplier: 1

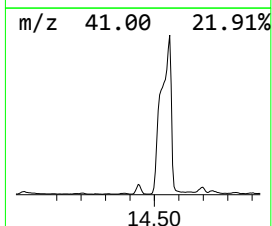
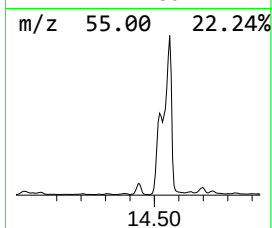
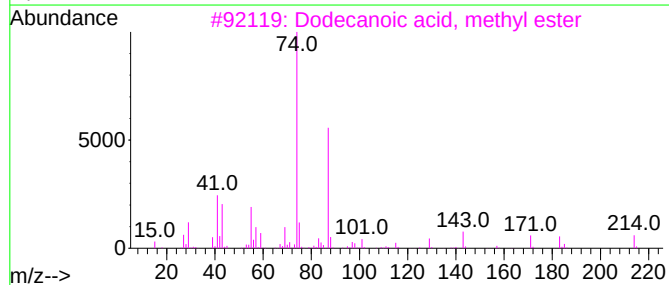
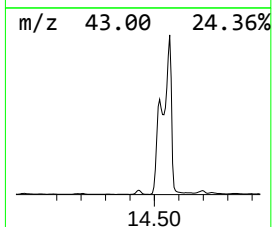
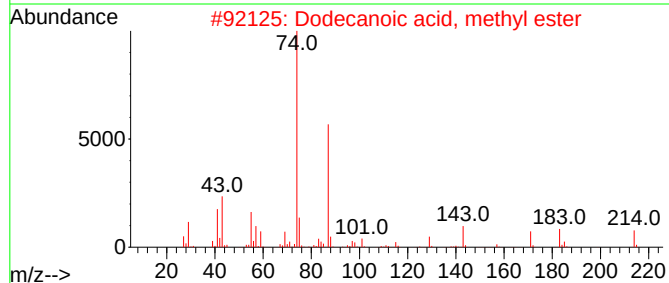
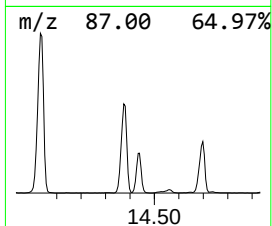
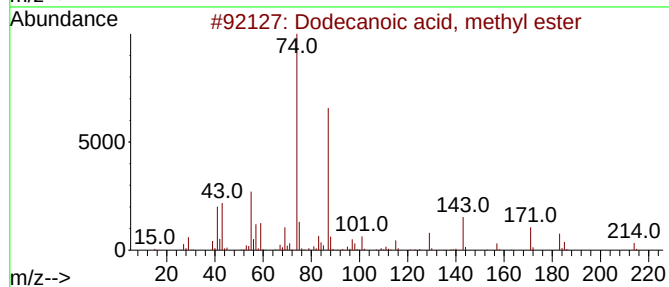
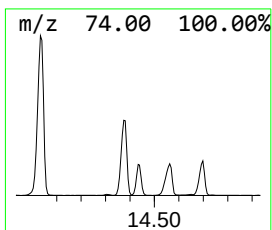
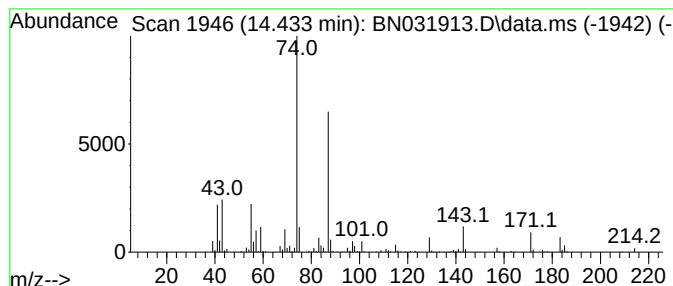
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Dodecanoic acid, methyl ester Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.433	2.11 ng/ul	379205	Acenaphthene-d10	14.310	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dodecanoic acid, methyl ester	214	C13H26O2	000111-82-0	96
2	Dodecanoic acid, methyl ester	214	C13H26O2	000111-82-0	91
3	Dodecanoic acid, methyl ester	214	C13H26O2	000111-82-0	91
4	Undecanoic acid, methyl ester	200	C12H24O2	001731-86-8	80
5	Nonanoic acid, methyl ester	172	C10H20O2	001731-84-6	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061224\
Data File : BN031913.D
Acq On : 11 Jun 2024 16:28
Operator : MA/JU
Sample : P2642-18
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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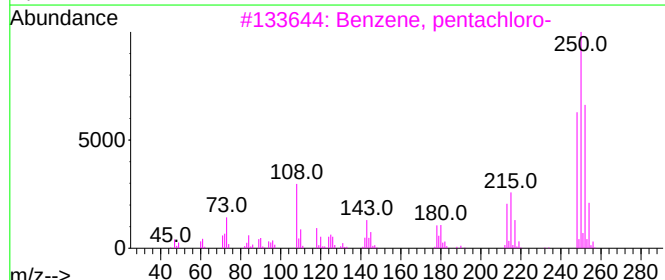
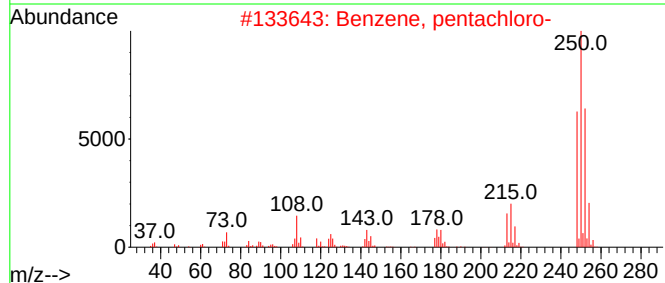
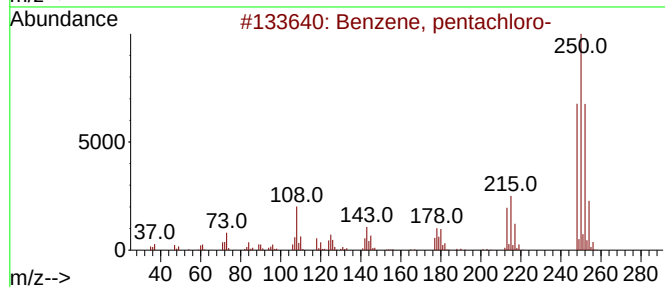
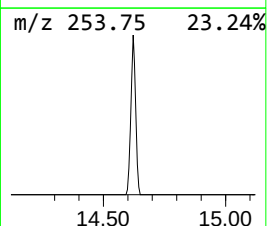
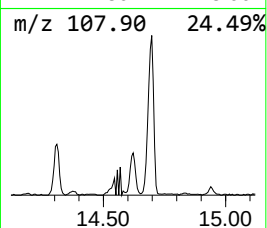
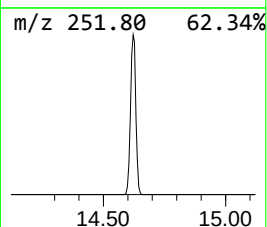
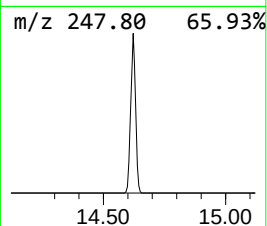
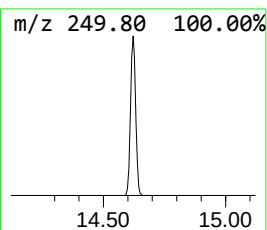
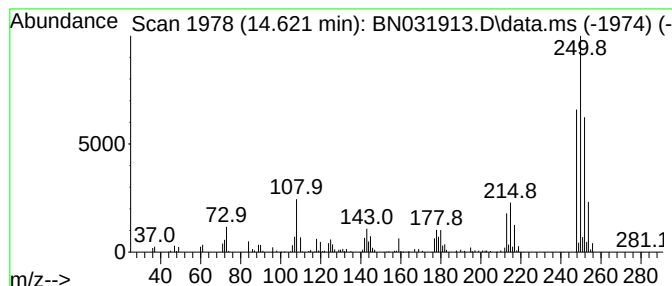
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Benzene, pentachloro- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.621	3.13 ng/ul	564272	Acenaphthene-d10	14.310

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, pentachloro-	248	C6HC15	000608-93-5	99
2			Benzene, pentachloro-	248	C6HC15	000608-93-5	99
3			Benzene, pentachloro-	248	C6HC15	000608-93-5	96
4			Benzene, pentachloro-	248	C6HC15	000608-93-5	95
5			Benzene, pentachloro-	248	C6HC15	000608-93-5	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061224\
Data File : BN031913.D
Acq On : 11 Jun 2024 16:28
Operator : MA/JU
Sample : P2642-18
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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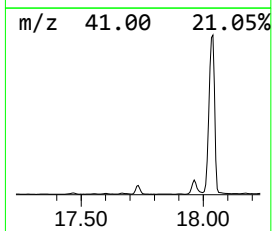
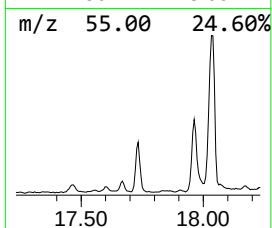
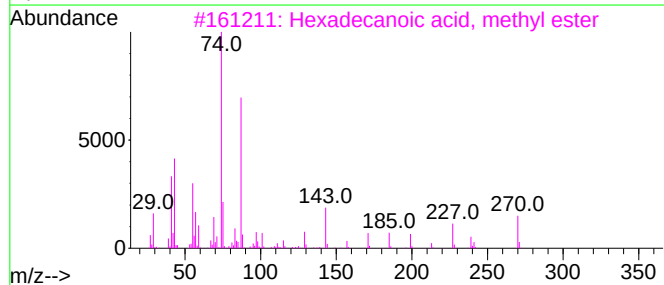
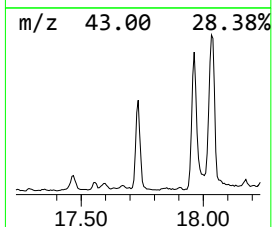
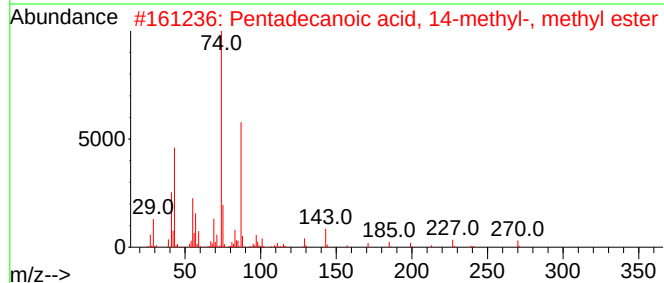
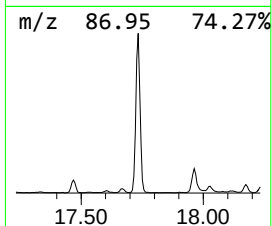
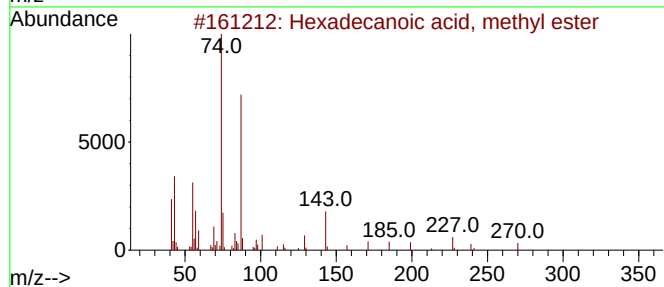
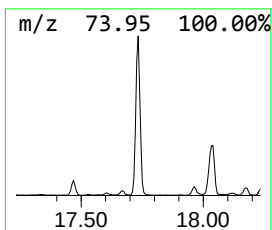
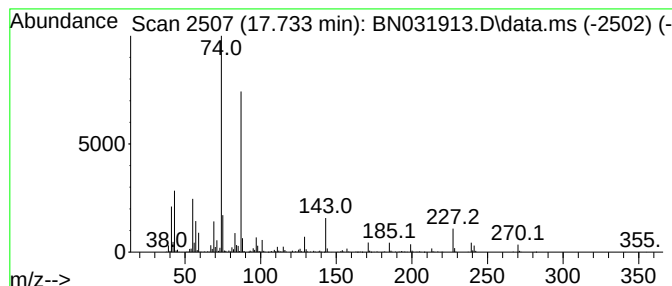
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Hexadecanoic acid, methyl e... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.733	4.68 ng/ul	1018430	Phenanthrene-d10	17.057

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	98
2			Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	97
3			Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	94
4			Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	91
5			Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061224\
Data File : BN031913.D
Acq On : 11 Jun 2024 16:28
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Sample : P2642-18
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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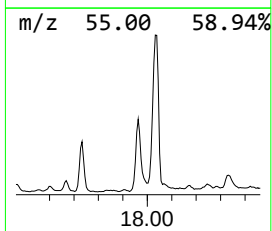
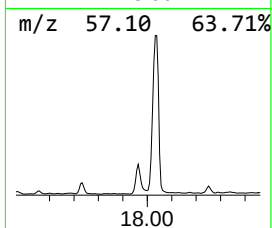
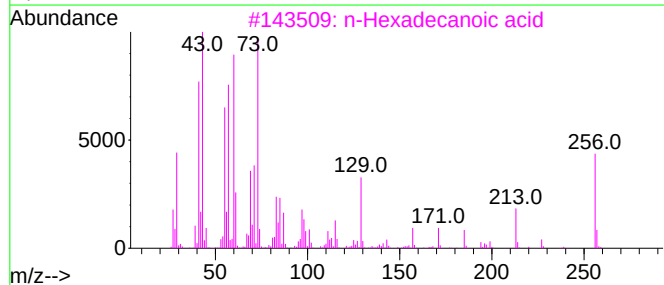
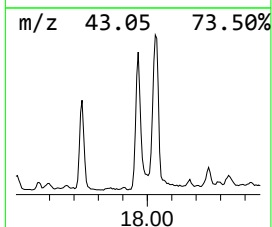
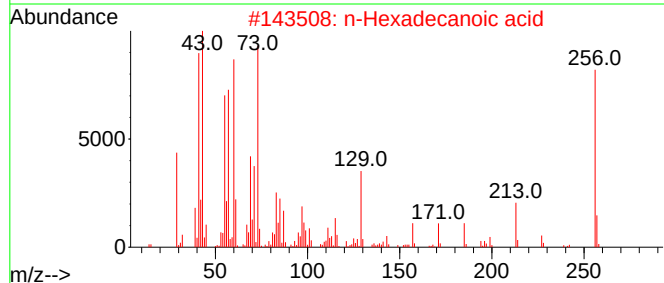
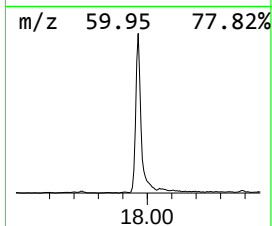
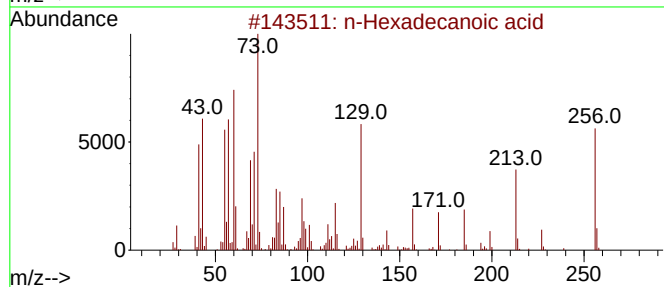
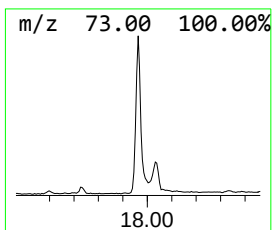
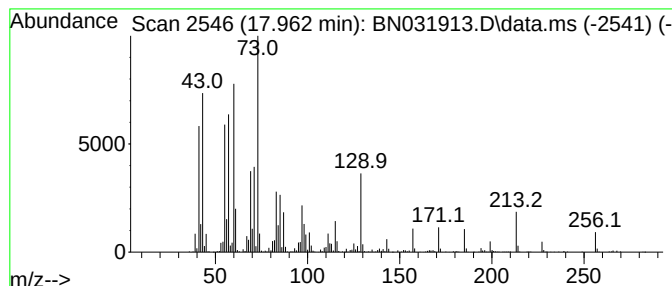
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.962	6.79 ng/ul	1478260	Phenanthrene-d10	17.057

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	97		
5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96		



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061224\
Data File : BN031913.D
Acq On : 11 Jun 2024 16:28
Operator : MA/JU
Sample : P2642-18
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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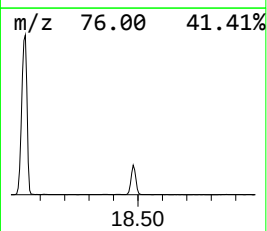
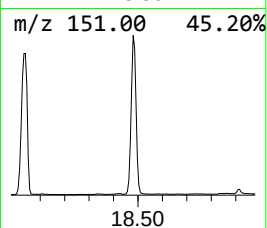
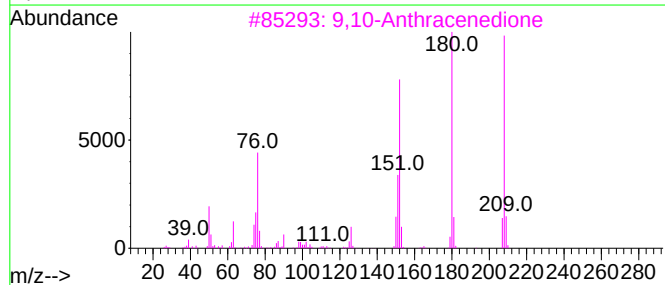
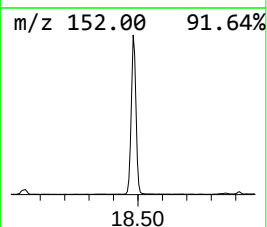
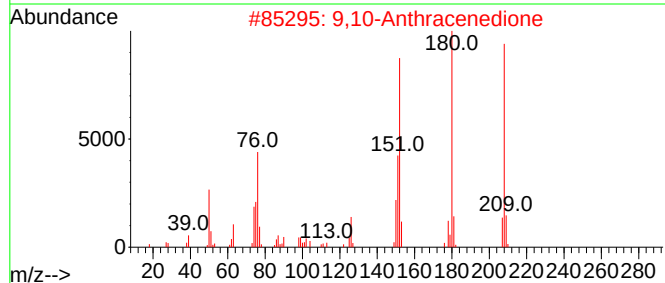
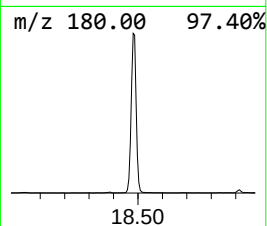
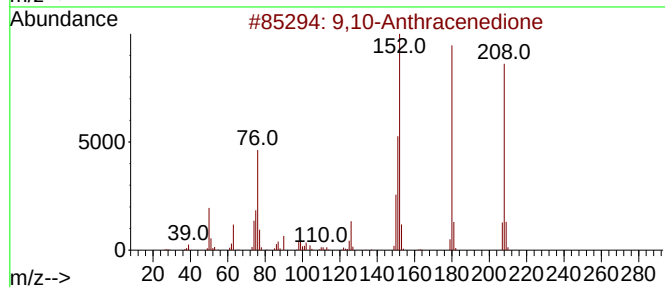
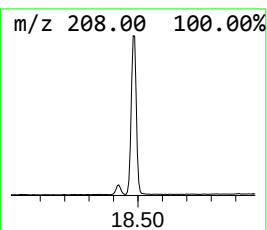
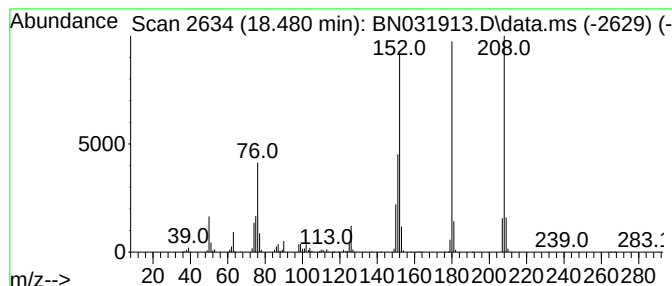
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 9,10-Anthracenedione Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.480	10.27 ng/ul	2235450	Phenanthrene-d10	17.057

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	98
4	9,10-Anthracenedione			208	C14H8O2	000084-65-1	95
5	9H-Fluoren-9-one			180	C13H8O	000486-25-9	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061224\
Data File : BN031913.D
Acq On : 11 Jun 2024 16:28
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Sample : P2642-18
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ALS Vial : 4 Sample Multiplier: 1

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ClientSampleId :
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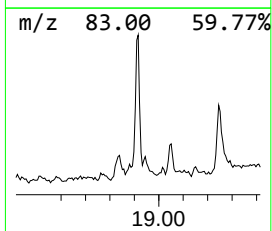
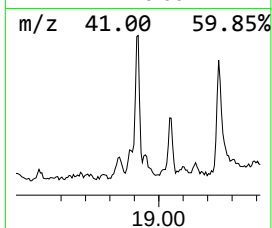
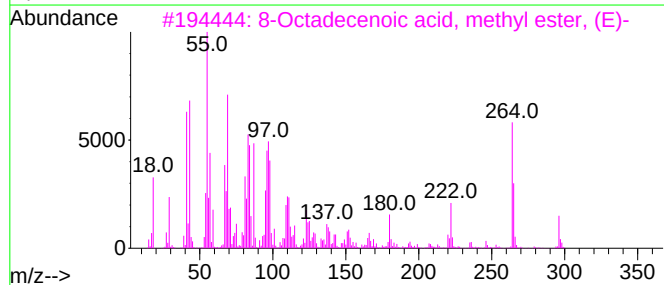
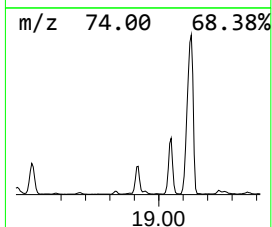
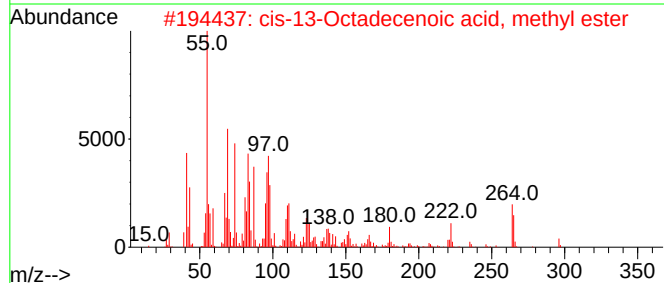
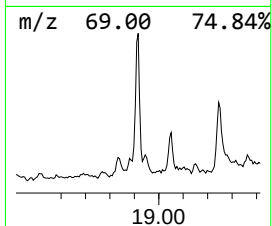
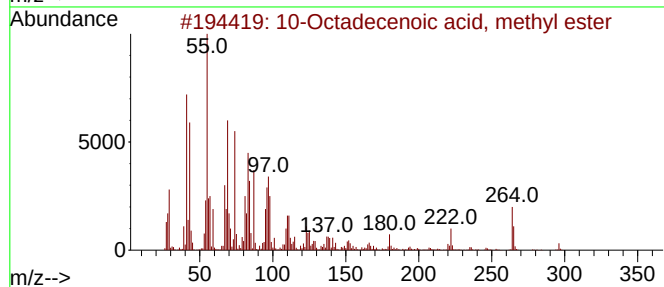
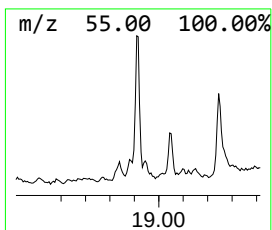
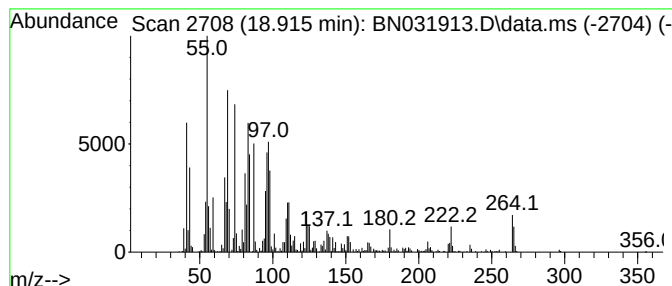
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 10-Octadecenoic acid, methy... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.915	4.06 ng/ul	883093	Phenanthrene-d10	17.057

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			10-Octadecenoic acid, methyl ester	296	C19H36O2	013481-95-3	99
2			cis-13-Octadecenoic acid, methyl...	296	C19H36O2	1010333-58-3	99
3			8-Octadecenoic acid, methyl este...	296	C19H36O2	026528-50-7	99
4			6-Octadecenoic acid, methyl ester	296	C19H36O2	052355-31-4	99
5			9-Octadecenoic acid, methyl ester	296	C19H36O2	002462-84-2	99



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061224\
Data File : BN031913.D
Acq On : 11 Jun 2024 16:28
Operator : MA/JU
Sample : P2642-18
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

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

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1-Phenoxypropan...	11.192	2.5	ng/ul	328593	2	10.445	2595050	20.0
unknown-02	12.939	3.2	ng/ul	580344	3	14.310	3602590	20.0
Dodecanoic acid...	14.433	2.1	ng/ul	379205	3	14.310	3602590	20.0
Benzene, pentac...	14.621	3.1	ng/ul	564272	3	14.310	3602590	20.0
Hexadecanoic ac...	17.733	4.7	ng/ul	1018430	4	17.057	4351800	20.0
n-Hexadecanoic ...	17.962	6.8	ng/ul	1478260	4	17.057	4351800	20.0
9,10-Anthracene...	18.480	10.3	ng/ul	2235450	4	17.057	4351800	20.0
10-Octadecenoic...	18.915	4.1	ng/ul	883093	4	17.057	4351800	20.0