

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102022\
 Data File : BP012250.D
 Acq On : 21 Oct 2022 17:49
 Operator : CG/JU
 Sample : N5064-12
 Misc :
 ALS Vial : 45 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0BH5

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_P\Methods\SFAM-EPA-BP101922.M
 Title : SVOA CALIBRATION

Signal : TIC: BP012250.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.258	43	46	54	rVB	68623	86640	1.25%	0.082%
2	3.781	131	135	144	rVB	237591	322365	4.66%	0.306%
3	3.905	151	156	162	rVB	96511	149746	2.16%	0.142%
4	3.999	167	172	181	rVB	56859	89458	1.29%	0.085%
5	4.922	324	329	339	rVB3	57257	132696	1.92%	0.126%
6	5.028	339	347	357	rBV3	32188	67584	0.98%	0.064%
7	5.452	415	419	427	rBV	39350	68337	0.99%	0.065%
8	5.822	476	482	490	rBV	62587	101210	1.46%	0.096%
9	6.681	623	628	639	rVB	30698	57997	0.84%	0.055%
10	6.993	670	681	694	rBV	1004197	1754694	25.35%	1.665%
11	7.158	702	709	729	rBV	1024314	1652943	23.88%	1.569%
12	7.352	735	742	752	rBV	1156510	1842506	26.62%	1.749%
13	7.822	811	822	831	rBV	1730744	2767310	39.97%	2.627%
14	8.540	937	944	953	rBV	992455	1740136	25.14%	1.652%
15	8.993	1013	1021	1028	rBV	1074491	1867487	26.98%	1.773%
16	9.381	1079	1087	1094	rBV2	38717	70868	1.02%	0.067%
17	9.710	1134	1143	1156	rBV	851208	1443297	20.85%	1.370%
18	10.252	1225	1235	1246	rBV	1226881	2163342	31.25%	2.053%
19	10.422	1258	1264	1273	rBV3	35168	84956	1.23%	0.081%
20	10.628	1288	1299	1304	rBV	2309832	4043582	58.41%	3.838%
21	10.675	1304	1307	1313	rVB	118327	187437	2.71%	0.178%
22	10.775	1318	1324	1342	rVB	319284	640176	9.25%	0.608%
23	12.287	1576	1581	1591	rVB	133519	219785	3.17%	0.209%
24	12.510	1613	1619	1628	rVB	131261	203837	2.94%	0.193%
25	13.287	1743	1751	1759	rBV2	58071	123427	1.78%	0.117%
26	13.363	1759	1764	1771	rBV	169582	263823	3.81%	0.250%
27	13.792	1831	1837	1842	rBV2	91926	167719	2.42%	0.159%
28	13.845	1842	1846	1848	rBV	65641	91145	1.32%	0.087%
29	13.887	1848	1853	1859	rVV	2194863	2990300	43.20%	2.838%
30	14.028	1873	1877	1880	rBV	49311	63836	0.92%	0.061%
31	14.163	1894	1900	1918	rBV2	2669492	4830581	69.78%	4.585%
32	14.475	1943	1953	1959	rBV	3247309	5073369	73.29%	4.815%
33	14.534	1959	1963	1971	rVB3	44317	97857	1.41%	0.093%
34	14.698	1983	1991	1997	rBV2	138876	298715	4.31%	0.284%
35	14.875	2013	2021	2030	rBV	125034	276316	3.99%	0.262%
36	15.087	2049	2057	2062	rBV3	42951	87966	1.27%	0.083%

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Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

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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_P\Methods\SFAM-EPA-BP101922.M
 Title : SVOA CALIBRATION

37	15.263	2080	2087	2090	rBV5	45500	90374	1.31%	0.086%
38	15.316	2090	2096	2099	rVV3	42033	75123	1.09%	0.071%
39	15.469	2115	2122	2127	rBV	3551986	4989591	72.08%	4.736%
40	15.522	2127	2131	2139	rVB3	135811	263247	3.80%	0.250%
41	15.598	2139	2144	2150	rBV	291893	400033	5.78%	0.380%
42	15.722	2161	2165	2175	rVB5	43132	98001	1.42%	0.093%
43	15.822	2176	2182	2187	rVV	86966	154183	2.23%	0.146%
44	15.963	2203	2206	2215	rVV	85517	173431	2.51%	0.165%
45	16.057	2215	2222	2228	rVB3	30390	68632	0.99%	0.065%
46	16.204	2240	2247	2253	rBV2	126377	235237	3.40%	0.223%
47	16.534	2295	2303	2308	rBV5	41542	99907	1.44%	0.095%
48	16.763	2337	2342	2346	rBV2	68794	117879	1.70%	0.112%
49	16.886	2358	2363	2371	rBV	212163	308332	4.45%	0.293%
50	16.957	2371	2375	2377	rBV	47973	68561	0.99%	0.065%
51	16.981	2377	2379	2383	rVV2	56514	65482	0.95%	0.062%
52	17.051	2386	2391	2399	rVB	110955	187293	2.71%	0.178%
53	17.233	2413	2422	2426	rBV	3520957	5182624	74.86%	4.919%
54	17.275	2426	2429	2434	rVV	2075853	2804498	40.51%	2.662%
55	17.333	2434	2439	2443	rVV	3039878	4546505	65.68%	4.315%
56	17.369	2443	2445	2450	rVB	427269	459499	6.64%	0.436%
57	17.645	2485	2492	2500	rBV	199124	351457	5.08%	0.334%
58	17.722	2502	2505	2508	rVV2	47065	63015	0.91%	0.060%
59	17.751	2508	2510	2516	rVV3	48629	71821	1.04%	0.068%
60	17.816	2517	2521	2528	rVB5	60183	104543	1.51%	0.099%
61	17.898	2531	2535	2540	rBV2	85426	130178	1.88%	0.124%
62	18.116	2565	2572	2574	rVV2	372059	771394	11.14%	0.732%
63	18.145	2574	2577	2582	rVV	442920	616455	8.90%	0.585%
64	18.228	2587	2591	2595	rVB	91616	121603	1.76%	0.115%
65	18.286	2595	2601	2605	rBV2	319345	603054	8.71%	0.572%
66	18.322	2605	2607	2614	rVB	226540	284249	4.11%	0.270%
67	18.398	2616	2620	2624	rBV4	96537	151887	2.19%	0.144%
68	18.480	2631	2634	2638	rBV4	50678	67906	0.98%	0.064%
69	18.586	2648	2652	2655	rBV	189347	225542	3.26%	0.214%
70	18.628	2655	2659	2670	rVB	373261	517076	7.47%	0.491%
71	18.822	2689	2692	2696	rBV2	56480	73140	1.06%	0.069%
72	18.869	2697	2700	2702	rVV	64882	73242	1.06%	0.070%
73	18.898	2703	2705	2710	rVB4	84778	127023	1.83%	0.121%
74	18.992	2716	2721	2722	rVV	170236	239161	3.45%	0.227%
75	19.033	2725	2728	2733	rVV4	302130	383891	5.55%	0.364%
76	19.080	2733	2736	2739	rVB	87232	97863	1.41%	0.093%

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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

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77	19.128	2740	2744	2753	rBV3	266273	523150	7.56%	0.497%
78	19.245	2756	2764	2772	rBV	2995060	4039668	58.35%	3.834%
79	19.351	2777	2782	2785	rBV3	265508	419382	6.06%	0.398%
80	19.386	2785	2788	2793	rVV2	405667	560936	8.10%	0.532%
81	19.439	2794	2797	2801	rVB2	94699	128581	1.86%	0.122%
82	19.569	2811	2819	2822	rBV	3642793	5531537	79.90%	5.250%
83	19.598	2822	2824	2832	rVV	2543097	3028371	43.75%	2.874%
84	19.669	2832	2836	2843	rVB2	183053	266044	3.84%	0.253%
85	19.775	2851	2854	2859	rVB	142334	190941	2.76%	0.181%
86	19.916	2873	2878	2884	rBV3	193371	473844	6.84%	0.450%
87	20.051	2896	2901	2902	rBV3	184600	284728	4.11%	0.270%
88	20.375	2953	2956	2959	rBV2	172035	181464	2.62%	0.172%
89	20.816	3027	3031	3036	rBV	556840	717019	10.36%	0.681%
90	20.974	3054	3058	3061	rBV2	244270	382287	5.52%	0.363%
91	21.027	3062	3067	3070	rBV2	471366	858950	12.41%	0.815%
92	21.104	3077	3080	3088	rVB	528768	704233	10.17%	0.668%
93	21.322	3111	3117	3121	rBV2	3907743	6922731	100.00%	6.571%
94	21.363	3121	3124	3134	rVB	1578190	2083484	30.10%	1.978%
95	21.539	3152	3154	3159	rVB2	233882	276886	4.00%	0.263%
96	22.998	3396	3402	3406	rBV	1606495	3655830	52.81%	3.470%
97	23.521	3486	3491	3495	rBV	1005482	1855673	26.81%	1.761%
98	23.574	3495	3500	3505	rVV2	2076149	4345530	62.77%	4.125%
99	23.621	3505	3508	3515	rVB	1291130	2301741	33.25%	2.185%
100	23.733	3521	3527	3533	rBV2	2152634	4036157	58.30%	3.831%

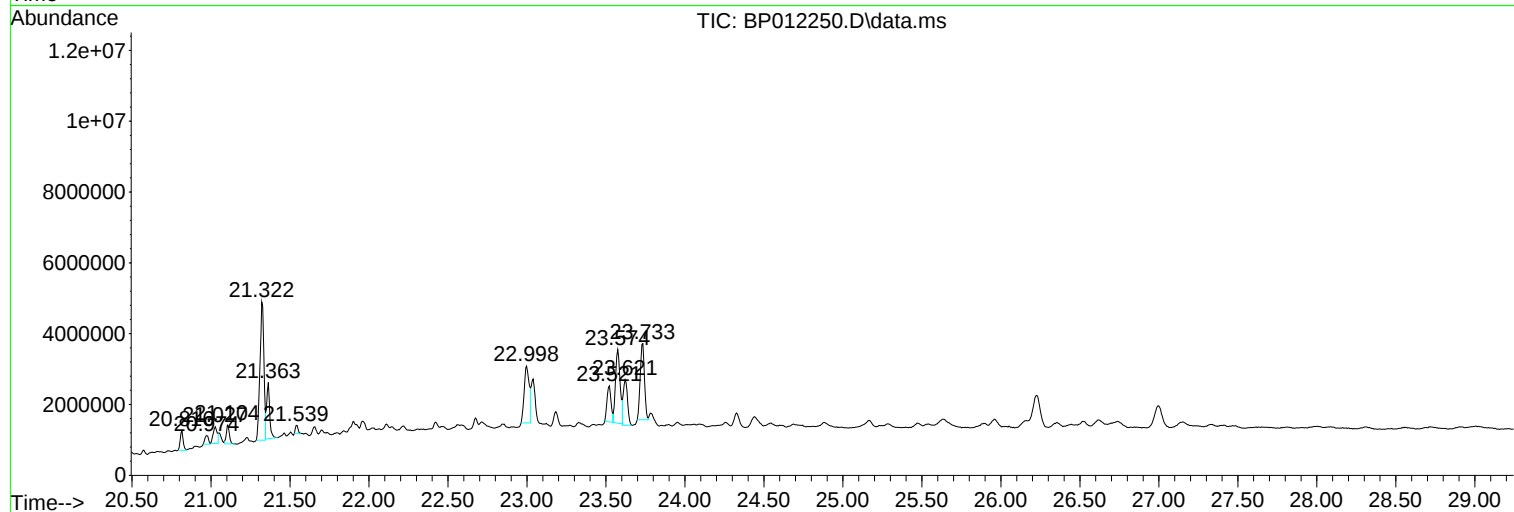
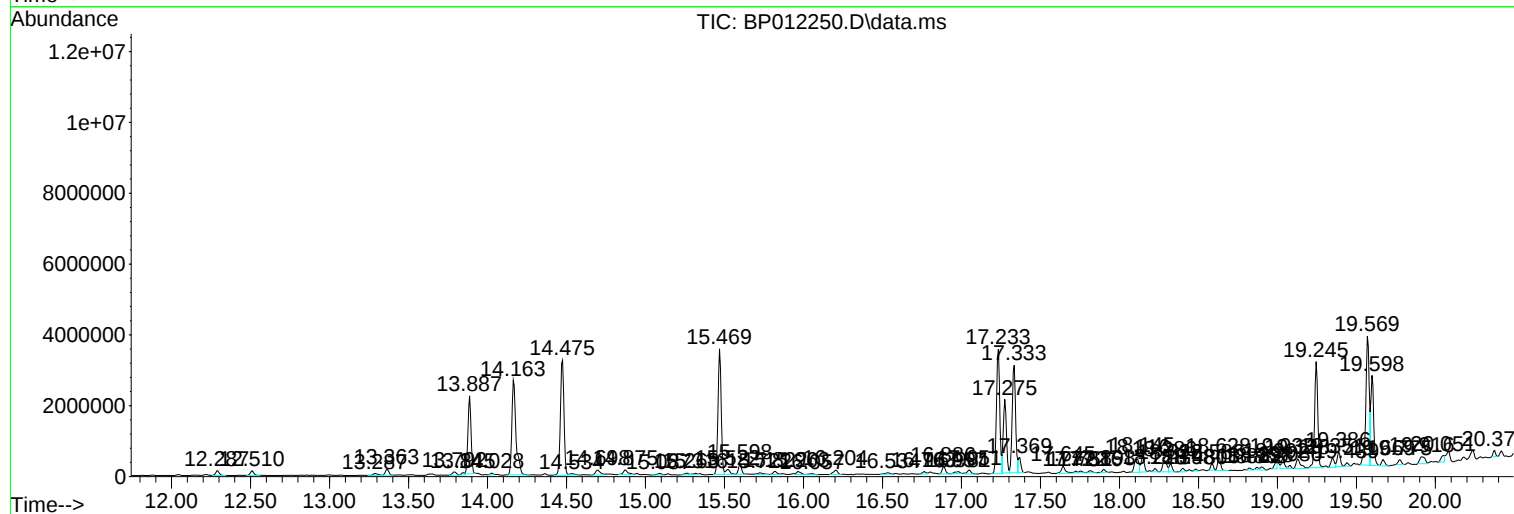
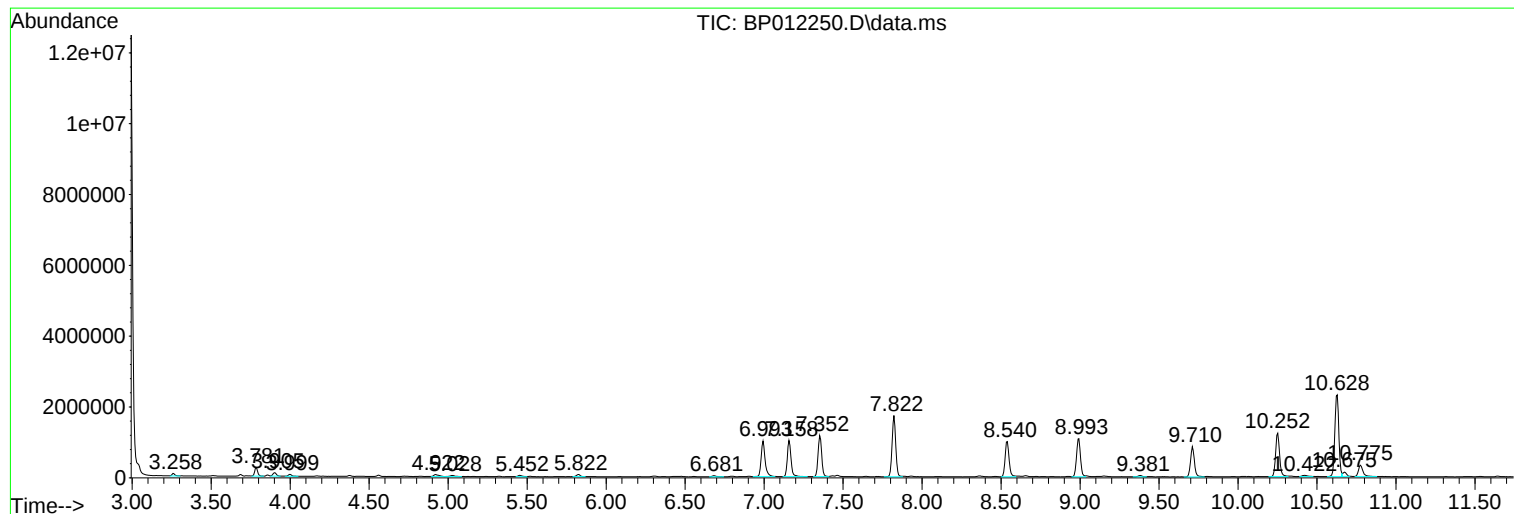
Sum of corrected areas: 105355542

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101922.M
Quant Title : SVOA CALIBRATION

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



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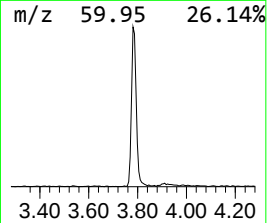
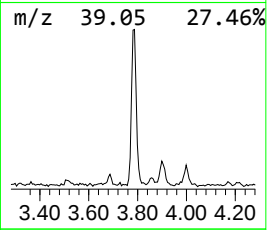
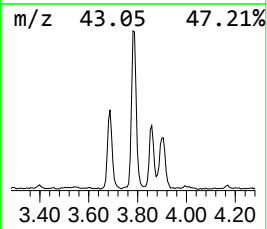
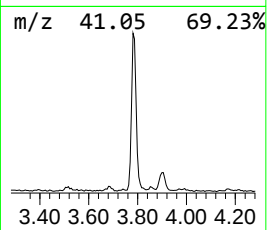
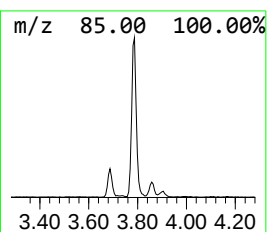
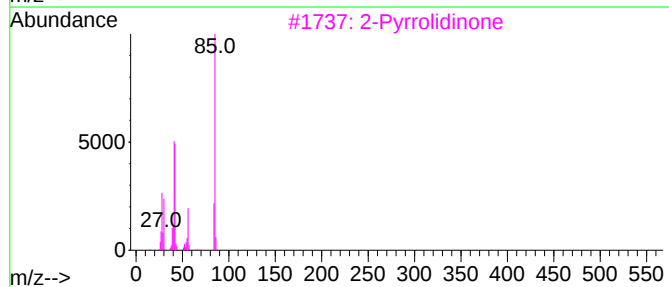
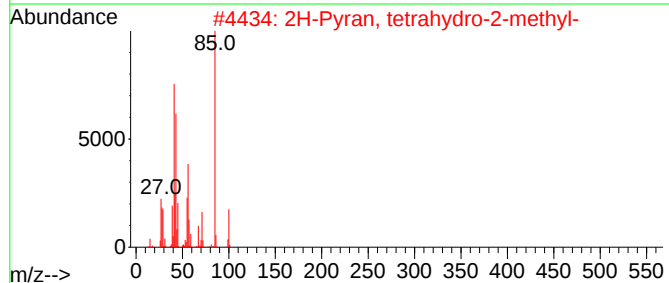
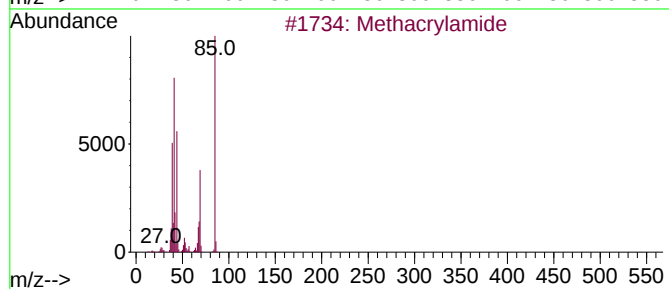
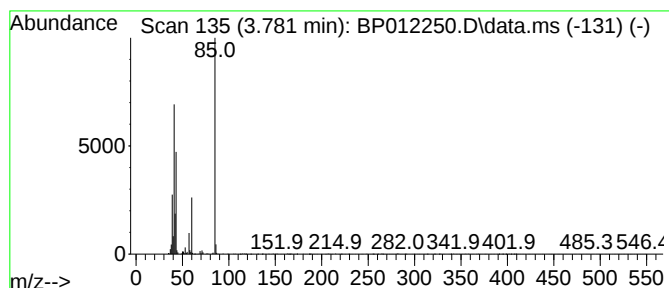
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101922.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 unknown-01 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.781	2.33 ng/ul	322365	1,4-Dichlorobenzene-d4	7.822

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methacrylamide	85	C4H7NO	000079-39-0	40
2			2H-Pyran, tetrahydro-2-methyl-	100	C6H12O	010141-72-7	38
3			2-Pyrrolidinone	85	C4H7NO	000616-45-5	9
4			Tetrahydropyran Z-10-dodecenoate	282	C17H30O3	1000130-96-5	9
5			2-Pyrrolidinone	85	C4H7NO	000616-45-5	9



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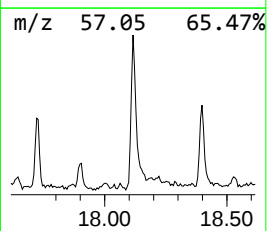
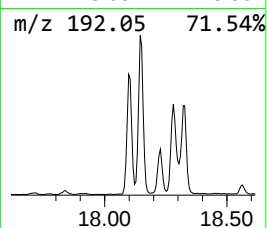
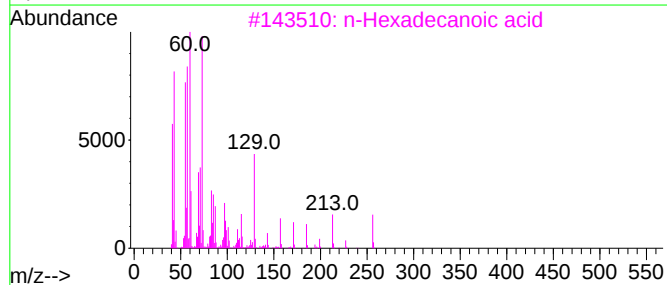
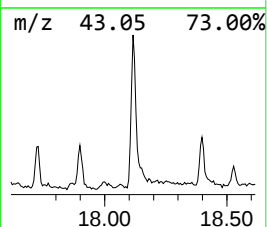
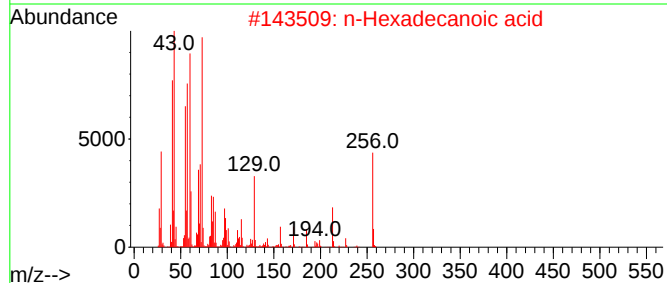
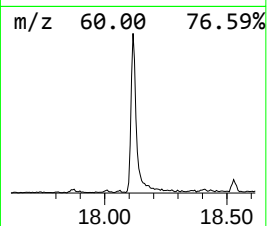
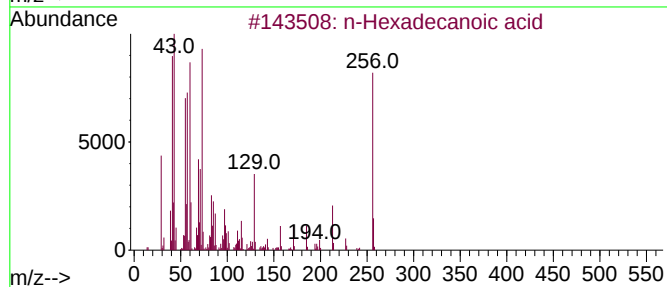
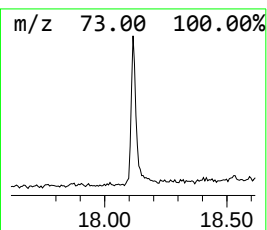
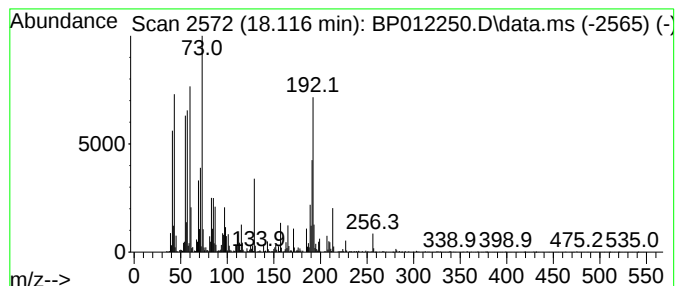
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101922.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.116	2.98 ng/ul	771394	Phenanthrene-d10	17.233

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
5			1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	90



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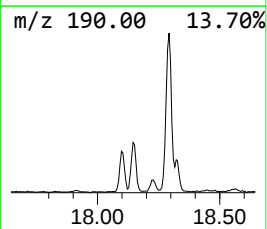
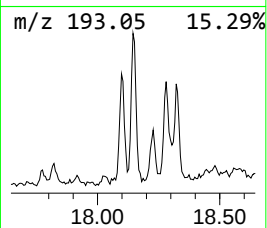
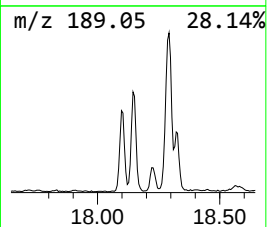
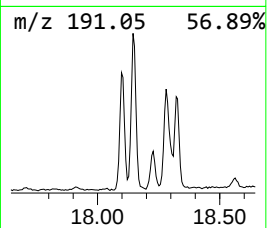
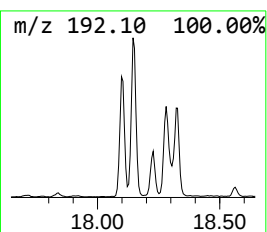
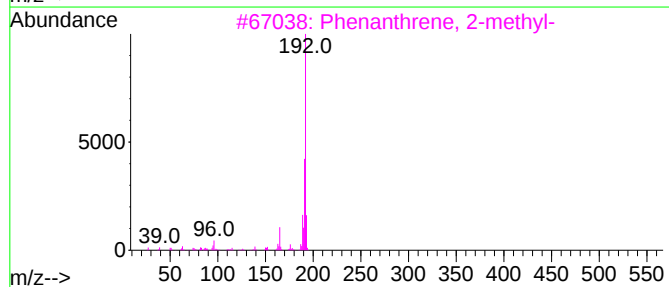
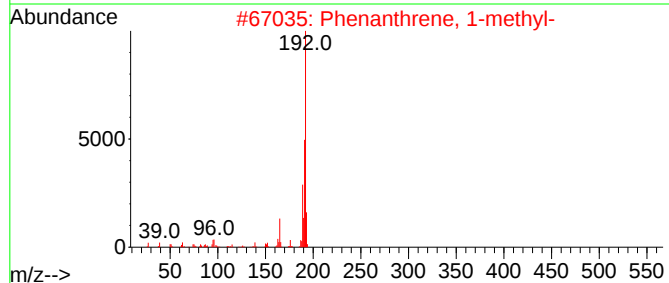
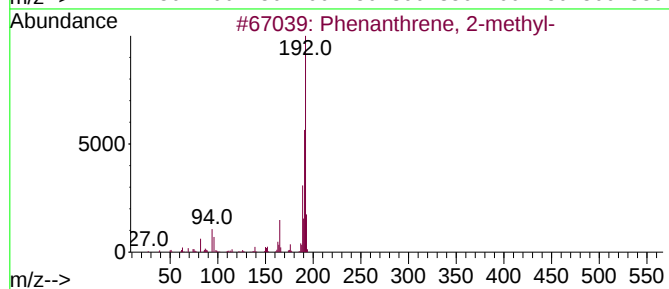
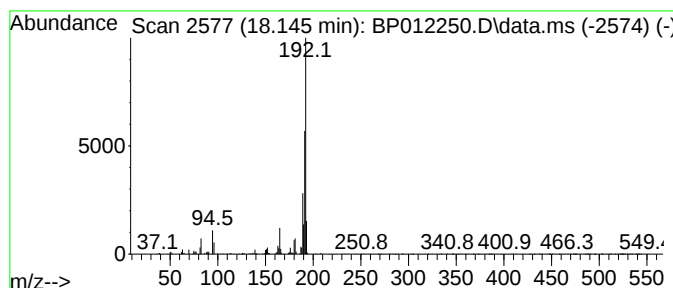
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Phenanthrene, 2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.145	2.38 ng/ul	616455	Phenanthrene-d10	17.233

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93
2			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	93
3			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93
4			1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	93
5			Anthracene, 2-methyl-	192	C15H12	000613-12-7	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102022\
Data File : BP012250.D
Acq On : 21 Oct 2022 17:49
Operator : CG/JU
Sample : N5064-12
Misc :
ALS Vial : 45 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0BH5

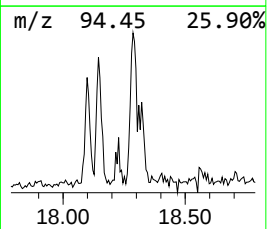
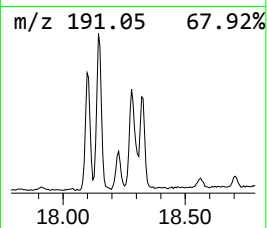
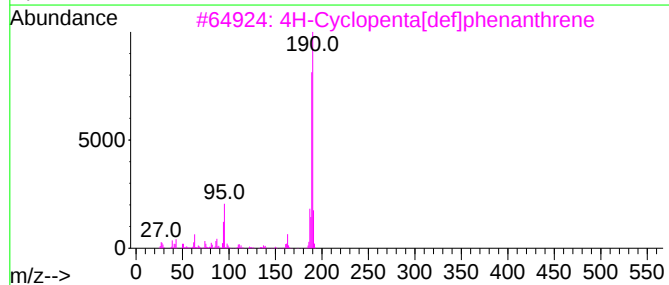
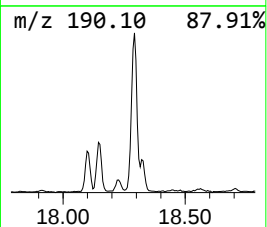
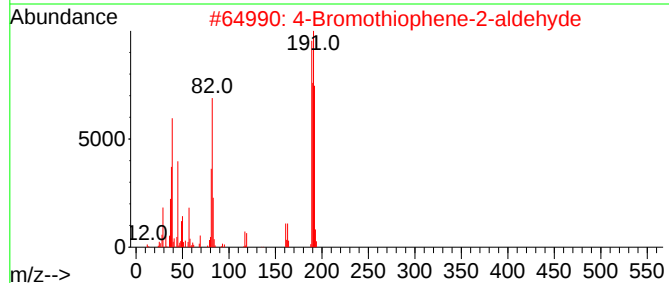
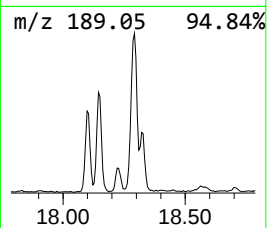
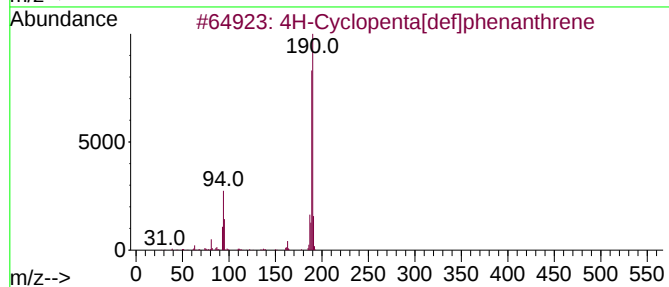
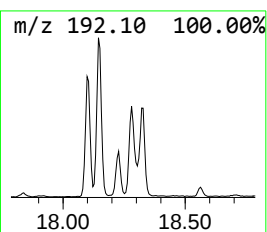
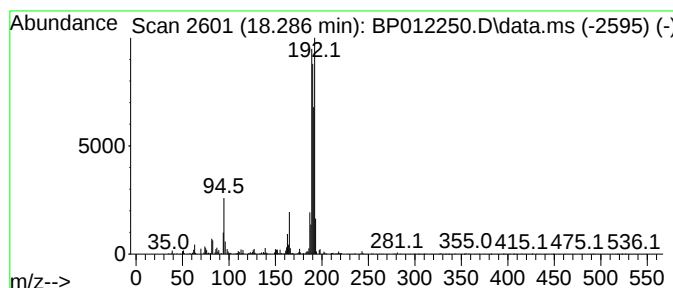
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 4H-Cyclopenta[def]phenanthrene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.286	2.33 ng/ul	603054	Phenanthrene-d10	17.233

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	60
2			4-Bromothiophene-2-aldehyde	190	C5H3BrOS	018791-75-8	58
3			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	46
4			Anthracene, 1-methyl-	192	C15H12	000610-48-0	46
5			Anthracene, 1-methyl-	192	C15H12	000610-48-0	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102022\
Data File : BP012250.D
Acq On : 21 Oct 2022 17:49
Operator : CG/JU
Sample : N5064-12
Misc :
ALS Vial : 45 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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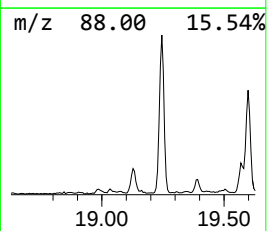
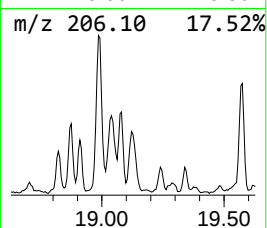
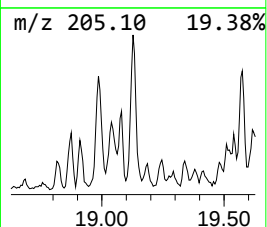
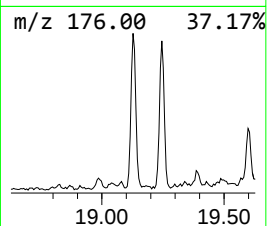
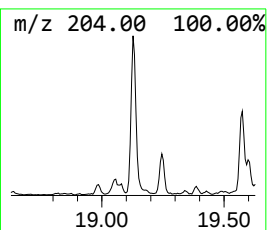
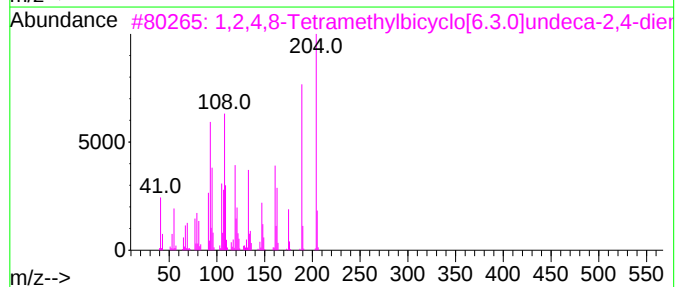
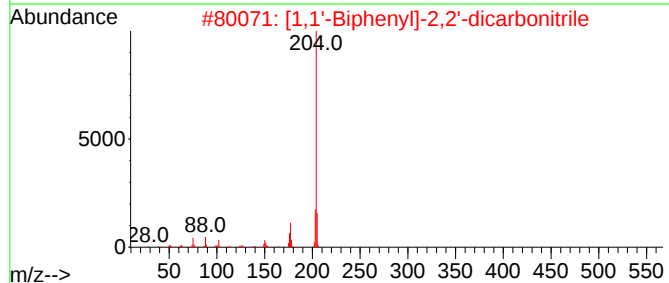
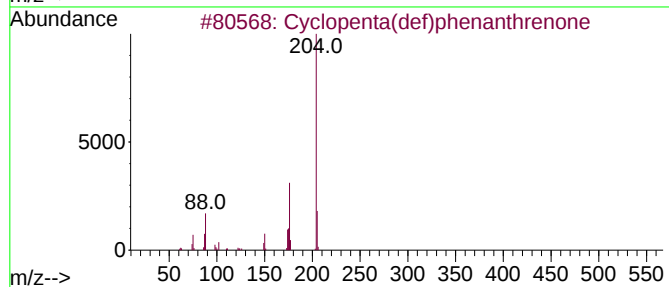
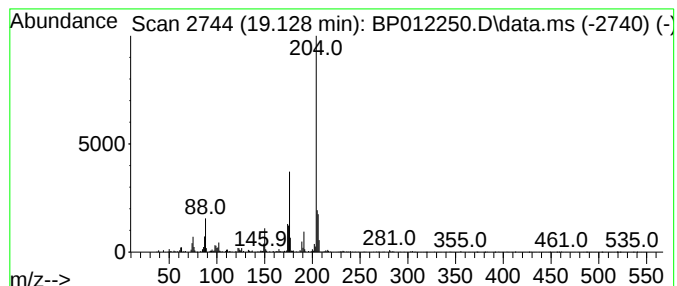
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Cyclopenta(def)phenanthrenone Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.128	2.02 ng/ul	523150	Phenanthrene-d10	17.233

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopenta(def)phenanthrenone	204	C15H8O	005737-13-3	96
2			[1,1'-Biphenyl]-2,2'-dicarbonitrile	204	C14H8N2	004341-02-0	50
3			1,2,4,8-Tetramethylbicyclo[6.3.0]undeca-2,4-dien-2-one	204	C15H24	137235-51-9	49
4			Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	49
5			9-Acridinecarbonitrile	204	C14H8N2	005326-19-2	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102022\
Data File : BP012250.D
Acq On : 21 Oct 2022 17:49
Operator : CG/JU
Sample : N5064-12
Misc :
ALS Vial : 45 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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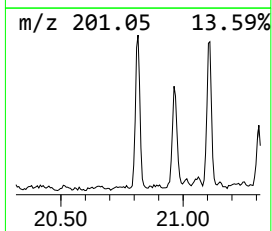
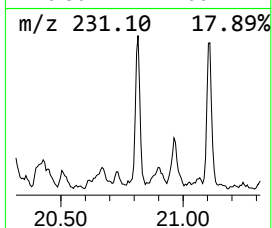
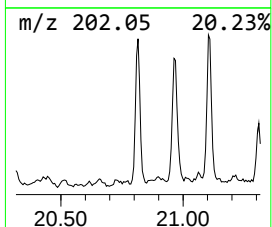
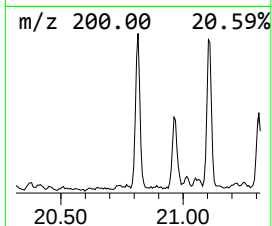
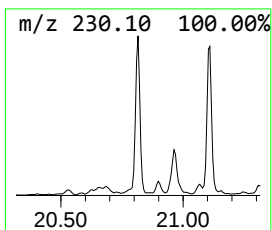
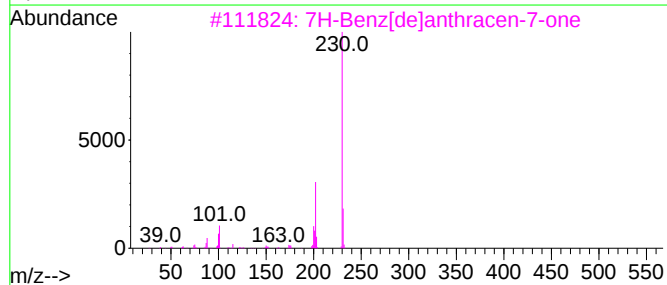
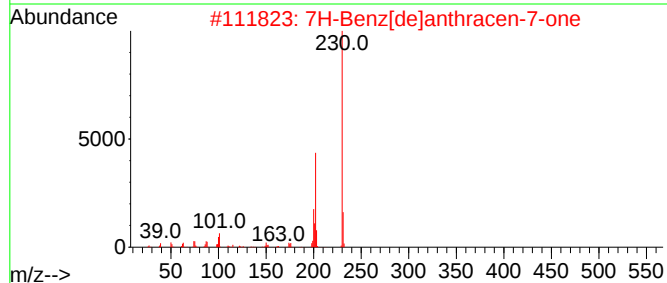
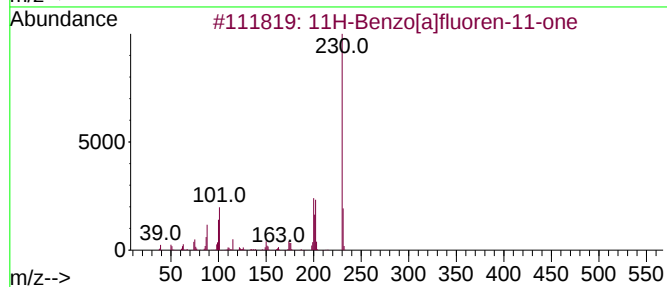
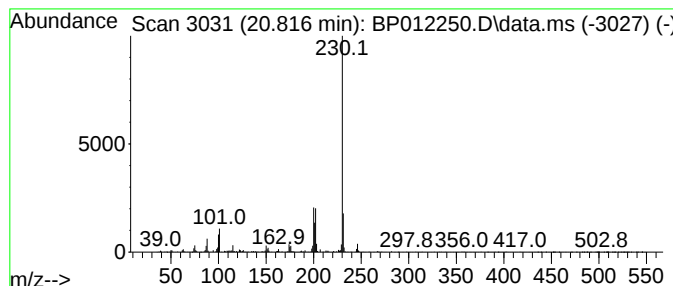
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 11H-Benzo[a]fluoren-11-one Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.816	2.07 ng/ul	717019	Chrysene-d12	21.327

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			11H-Benzo[a]fluoren-11-one	230	C17H10O	000479-79-8	97
2			7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	90
3			7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	89
4			7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	87
5			7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102022\
Data File : BP012250.D
Acq On : 21 Oct 2022 17:49
Operator : CG/JU
Sample : N5064-12
Misc :
ALS Vial : 45 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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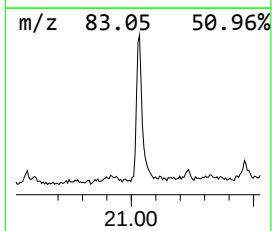
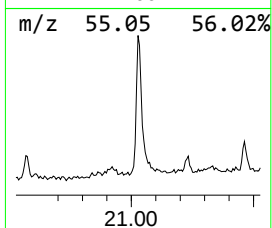
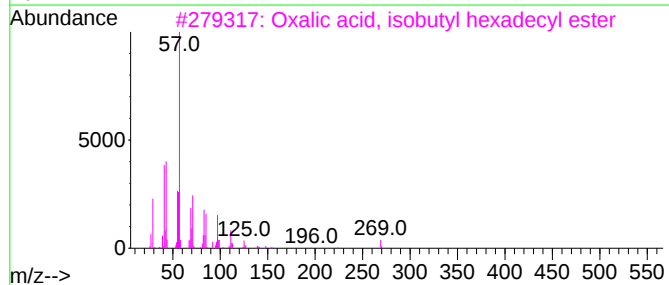
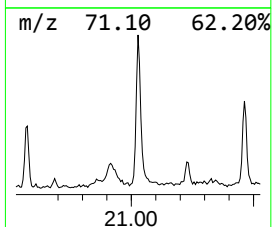
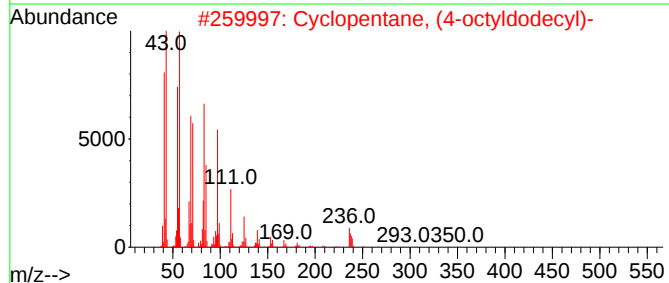
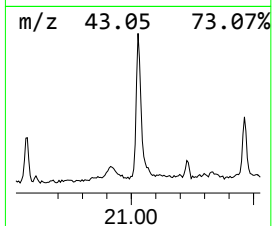
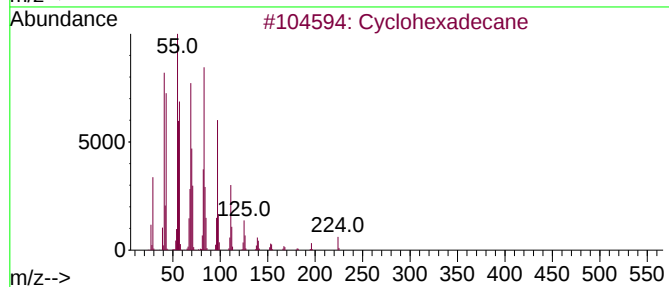
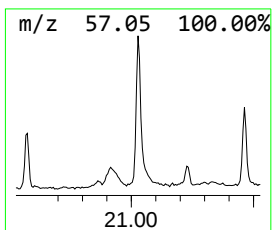
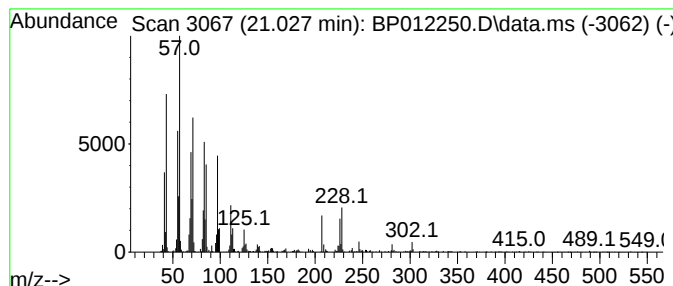
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Cyclic21.027 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.027	2.48 ng/ul	858950	Chrysene-d12	21.327

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexadecane	224	C16H32	000295-65-8	92
2			Cyclopentane, (4-octyldodecyl)-	350	C25H50	005638-09-5	90
3			Oxalic acid, isobutyl hexadecyl ...	370	C22H42O4	1000309-38-1	74
4			1-Decanol, 2-hexyl-	242	C16H34O	002425-77-6	70
5			Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102022\
Data File : BP012250.D
Acq On : 21 Oct 2022 17:49
Operator : CG/JU
Sample : N5064-12
Misc :
ALS Vial : 45 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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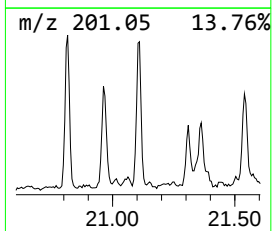
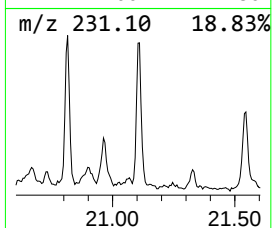
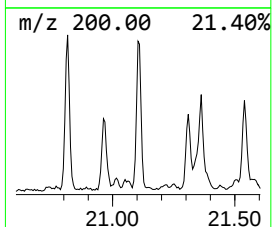
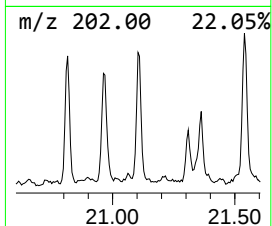
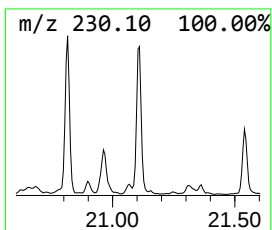
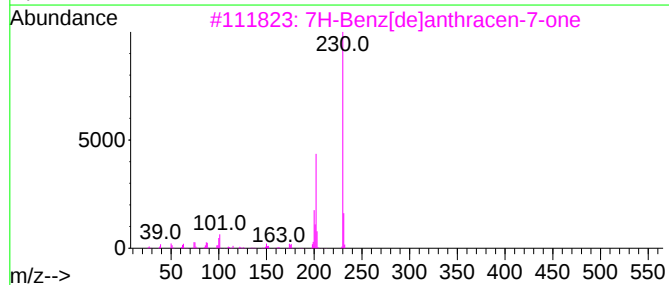
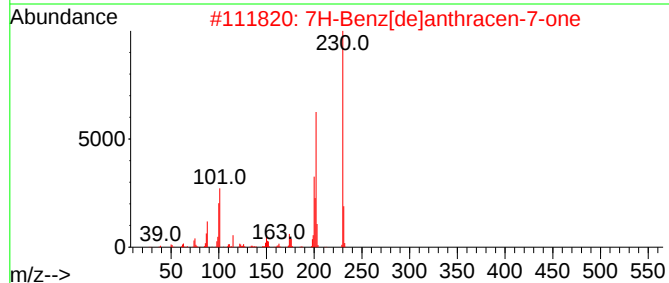
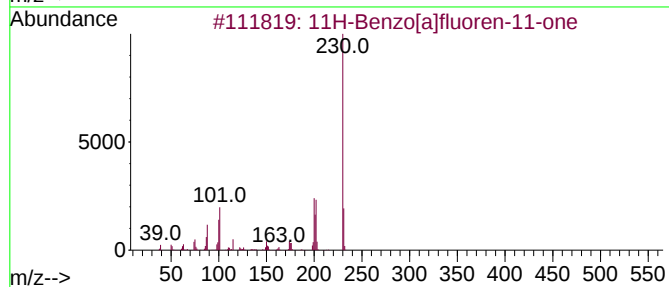
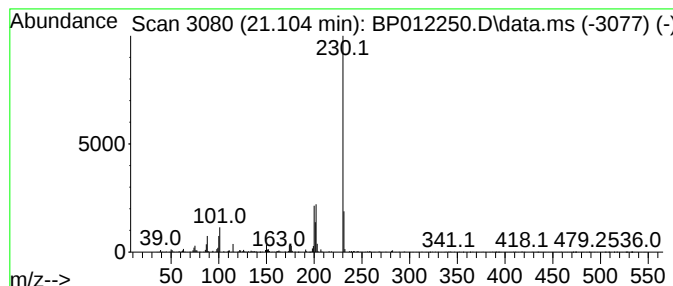
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 7H-Benz[de]anthracen-7-one Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.104	2.03 ng/ul	704233	Chrysene-d12	21.327

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	11H-Benzo[a]fluoren-11-one			230	C17H10O	000479-79-8	95
2	7H-Benz[de]anthracen-7-one			230	C17H10O	000082-05-3	90
3	7H-Benz[de]anthracen-7-one			230	C17H10O	000082-05-3	90
4	7H-Benz[de]anthracen-7-one			230	C17H10O	000082-05-3	90
5	7H-Benz[de]anthracen-7-one			230	C17H10O	000082-05-3	74



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102022\
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Acq On : 21 Oct 2022 17:49
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Sample : N5064-12
Misc :
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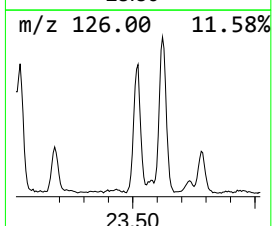
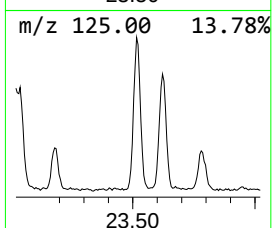
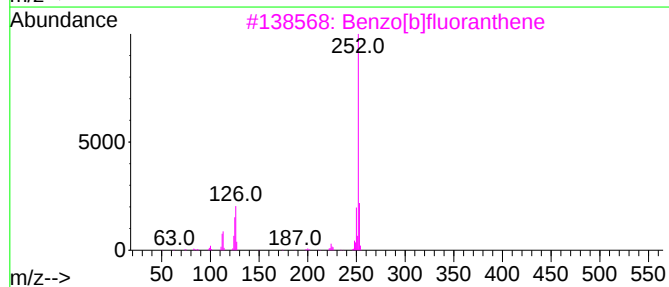
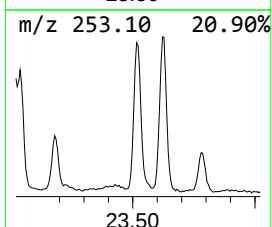
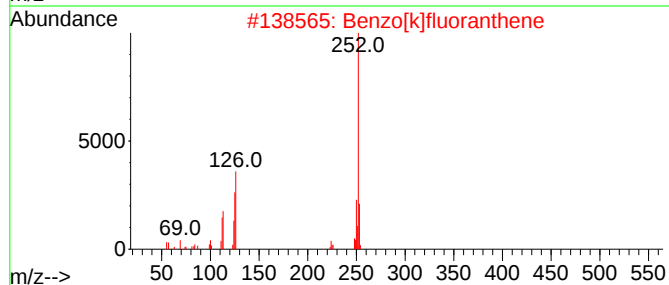
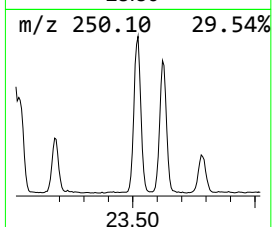
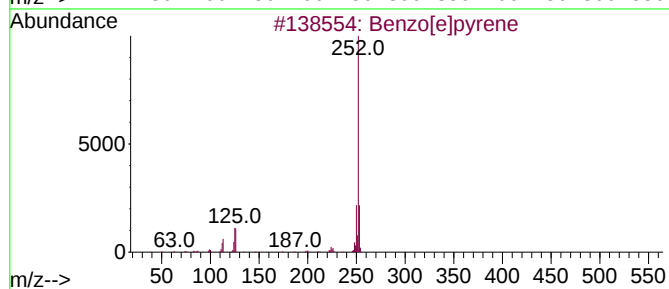
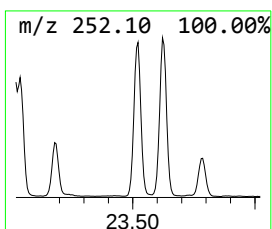
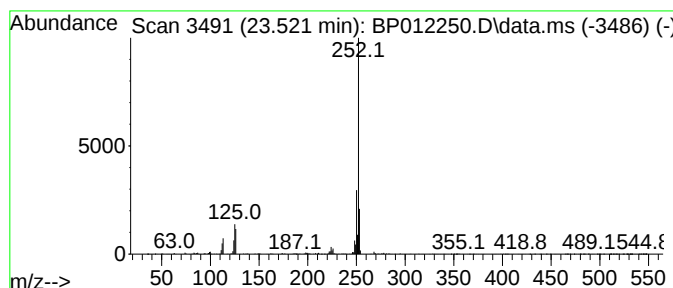
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Benzo[e]pyrene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.521	9.20 ng/ul	1855670	Perylene-d12	23.733

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[e]pyrene	252	C20H12	000192-97-2	98
2			Benzo[k]fluoranthene	252	C20H12	000207-08-9	95
3			Benzo[b]fluoranthene	252	C20H12	000205-99-2	95
4			Benzo[b]fluoranthene	252	C20H12	000205-99-2	95
5			Benzo[a]pyrene	252	C20H12	000050-32-8	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102022\
Data File : BP012250.D
Acq On : 21 Oct 2022 17:49
Operator : CG/JU
Sample : N5064-12
Misc :
ALS Vial : 45 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101922.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
unknown-01	3.781	2.3	ng/u1	322365	1	7.822	2767310	20.0
n-Hexadecanoic ...	18.116	3.0	ng/u1	771394	4	17.233	5182620	20.0
Phenanthrene, 2...	18.145	2.4	ng/u1	616455	4	17.233	5182620	20.0
4H-Cyclopenta[d...	18.286	2.3	ng/u1	603054	4	17.233	5182620	20.0
Cyclopenta(def)...	19.128	2.0	ng/u1	523150	4	17.233	5182620	20.0
11H-Benzo[a]flu...	20.816	2.1	ng/u1	717019	5	21.327	6922730	20.0
(DEL) Alkane: C...	21.027	2.5	ng/u1	858950	5	21.327	6922730	20.0
7H-Benz[de]anth...	21.104	2.0	ng/u1	704233	5	21.327	6922730	20.0
Benzo[e]pyrene	23.521	9.2	ng/u1	1855670	6	23.733	4036160	20.0