

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110923\
 Data File : BP018015.D
 Acq On : 10 Nov 2023 06:51
 Operator : MA/JU
 Sample : 05091-01
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCK10

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-BP110923.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP018015.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.246	24	27	33	rVB	12135	17397	0.62%	0.055%
2	3.711	102	106	115	rVB2	10005	22007	0.79%	0.070%
3	3.793	116	120	124	rVV2	17465	23556	0.84%	0.075%
4	3.858	127	131	138	rVB	17552	28458	1.02%	0.090%
5	3.975	149	151	156	rVB5	6381	9038	0.32%	0.029%
6	4.564	248	251	259	rVB8	5766	10596	0.38%	0.034%
7	4.911	306	310	317	rBV	20935	30502	1.09%	0.097%
8	5.822	460	465	472	rBV5	6678	12543	0.45%	0.040%
9	7.034	663	671	689	rBV	153237	378181	13.50%	1.199%
10	7.216	696	702	716	rBV	120396	208723	7.45%	0.662%
11	7.387	724	731	742	rBV	176744	302997	10.82%	0.961%
12	7.863	805	812	823	rBV	319734	517485	18.48%	1.641%
13	8.399	897	903	914	rVB2	11874	26493	0.95%	0.084%
14	8.593	928	936	951	rBV	149462	308724	11.02%	0.979%
15	8.722	951	958	969	rVV	224239	372368	13.30%	1.181%
16	9.069	1008	1017	1027	rBV	166461	292300	10.44%	0.927%
17	9.787	1132	1139	1155	rBV	141362	270016	9.64%	0.856%
18	9.963	1162	1169	1182	rBV2	24267	75857	2.71%	0.241%
19	10.157	1197	1202	1214	rVB3	20461	45778	1.63%	0.145%
20	10.322	1223	1230	1251	rVB	169500	385449	13.76%	1.222%
21	10.687	1284	1292	1297	rBV	410430	701208	25.04%	2.224%
22	10.734	1297	1300	1309	rVB	22084	39633	1.42%	0.126%
23	10.875	1317	1324	1338	rBV	50780	132738	4.74%	0.421%
24	11.275	1385	1392	1397	rBV8	8161	23453	0.84%	0.074%
25	11.740	1463	1471	1478	rBV8	4655	11957	0.43%	0.038%
26	12.134	1534	1538	1551	rVB4	6197	17799	0.64%	0.056%
27	12.263	1554	1560	1568	rVB3	8887	27340	0.98%	0.087%
28	12.351	1570	1575	1579	rBV	12316	19798	0.71%	0.063%
29	12.581	1608	1614	1621	rVB7	5128	10459	0.37%	0.033%
30	12.845	1655	1659	1665	rVB9	4480	8027	0.29%	0.025%
31	13.375	1743	1749	1755	rBV2	24731	42038	1.50%	0.133%
32	13.492	1763	1769	1773	rBV8	6407	13021	0.46%	0.041%
33	13.957	1841	1848	1861	rVB	432741	615716	21.98%	1.953%
34	14.087	1866	1870	1878	rVB8	5232	9703	0.35%	0.031%
35	14.234	1887	1895	1913	rBV2	445603	770668	27.52%	2.444%
36	14.363	1913	1917	1921	rVB7	5696	8174	0.29%	0.026%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110923.MA.M
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37	14.539	1940	1947	1954	rVV	630692	953047	34.03%	3.022%
38	14.598	1954	1957	1965	rVB	39439	64914	2.32%	0.206%
39	14.810	1987	1993	2010	rBV2	62885	206517	7.37%	0.655%
40	14.951	2010	2017	2020	rVV2	19042	36733	1.31%	0.116%
41	15.145	2045	2050	2054	rBV7	6372	11609	0.41%	0.037%
42	15.316	2075	2079	2083	rVV2	20749	25890	0.92%	0.082%
43	15.363	2083	2087	2093	rVB3	10152	14371	0.51%	0.046%
44	15.539	2108	2117	2123	rBV	609225	880637	31.44%	2.793%
45	15.598	2123	2127	2135	rVB	57384	84737	3.03%	0.269%
46	15.692	2138	2143	2151	rBV3	160407	291900	10.42%	0.926%
47	15.886	2172	2176	2179	rBV5	10235	14007	0.50%	0.044%
48	16.033	2196	2201	2204	rBV6	10160	18960	0.68%	0.060%
49	16.133	2214	2218	2222	rVB7	7554	11216	0.40%	0.036%
50	16.192	2223	2228	2232	rBV2	52450	84412	3.01%	0.268%
51	16.310	2243	2248	2254	rBV2	22987	34377	1.23%	0.109%
52	16.657	2302	2307	2311	rBV8	9798	23398	0.84%	0.074%
53	16.957	2354	2358	2360	rBV5	25284	39818	1.42%	0.126%
54	16.992	2360	2364	2367	rVV2	73359	102182	3.65%	0.324%
55	17.033	2367	2371	2376	rVB	55852	81132	2.90%	0.257%
56	17.139	2385	2389	2392	rBV2	18761	24897	0.89%	0.079%
57	17.233	2401	2405	2409	rBV5	14850	22977	0.82%	0.073%
58	17.316	2413	2419	2423	rVV	856766	1204707	43.01%	3.821%
59	17.357	2423	2426	2431	rVV	220306	309764	11.06%	0.982%
60	17.416	2431	2436	2439	rVV2	711823	962302	34.36%	3.052%
61	17.451	2439	2442	2452	rVV	451426	681922	24.35%	2.163%
62	17.533	2452	2456	2463	rVB	30201	51964	1.86%	0.165%
63	17.745	2487	2492	2499	rVB2	123622	217543	7.77%	0.690%
64	17.898	2515	2518	2522	rBV5	15873	24890	0.89%	0.079%
65	18.163	2559	2563	2571	rBV2	243638	394575	14.09%	1.251%
66	18.233	2572	2575	2584	rVV3	50491	98157	3.50%	0.311%
67	18.310	2585	2588	2594	rVV	52452	73455	2.62%	0.233%
68	18.375	2594	2599	2603	rVV2	139919	228270	8.15%	0.724%
69	18.416	2603	2606	2615	rVB3	83511	124280	4.44%	0.394%
70	18.675	2646	2650	2654	rVB	50328	66065	2.36%	0.210%
71	18.739	2657	2661	2664	rVV2	89839	135939	4.85%	0.431%
72	18.769	2664	2666	2671	rVB2	76233	85798	3.06%	0.272%
73	18.992	2700	2704	2707	rBV2	76255	104055	3.72%	0.330%
74	19.033	2708	2711	2714	rVV2	58865	81603	2.91%	0.259%
75	19.069	2714	2717	2721	rVB	43632	57596	2.06%	0.183%
76	19.233	2741	2745	2757	rVB	302359	545297	19.47%	1.729%

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77	19.339	2758	2763	2769	rVB	1196114	1541939	55.06%	4.890%
78	19.404	2770	2774	2779	rBV2	148269	220412	7.87%	0.699%
79	19.669	2812	2819	2821	rBV2	1200723	1705689	60.90%	5.409%
80	19.698	2821	2824	2831	rVV	1303581	1636530	58.43%	5.190%
81	19.757	2831	2834	2838	rVB2	47189	69400	2.48%	0.220%
82	19.869	2850	2853	2857	rBV	107650	128204	4.58%	0.407%
83	20.004	2871	2876	2884	rVB2	98719	233780	8.35%	0.741%
84	20.151	2894	2901	2903	rBV5	122276	257182	9.18%	0.816%
85	20.180	2903	2906	2910	rVB	192051	249352	8.90%	0.791%
86	20.274	2919	2922	2925	rBV2	161730	193381	6.90%	0.613%
87	20.469	2952	2955	2960	rBV	125445	168992	6.03%	0.536%
88	20.516	2961	2963	2970	rVB3	76207	108176	3.86%	0.343%
89	21.080	3056	3059	3062	rBV2	204191	273949	9.78%	0.869%
90	21.121	3063	3066	3069	rVV2	195875	265208	9.47%	0.841%
91	21.157	3069	3072	3078	rVB2	164523	233099	8.32%	0.739%
92	21.439	3110	3120	3124	rBV2	1392181	2800713	100.00%	8.882%
93	21.480	3124	3127	3132	rVB	748907	947802	33.84%	3.006%
94	21.998	3213	3215	3218	rBV4	82352	107714	3.85%	0.342%
95	22.657	3324	3327	3331	rVB	175548	229707	8.20%	0.728%
96	23.174	3409	3415	3421	rBV	701512	1735319	61.96%	5.503%
97	23.727	3504	3509	3514	rBV	430127	854837	30.52%	2.711%
98	23.786	3514	3519	3524	rVV2	646503	1270723	45.37%	4.030%
99	23.839	3524	3528	3534	rVB2	331729	625350	22.33%	1.983%
100	23.951	3541	3547	3553	rBV	706499	1408710	50.30%	4.468%

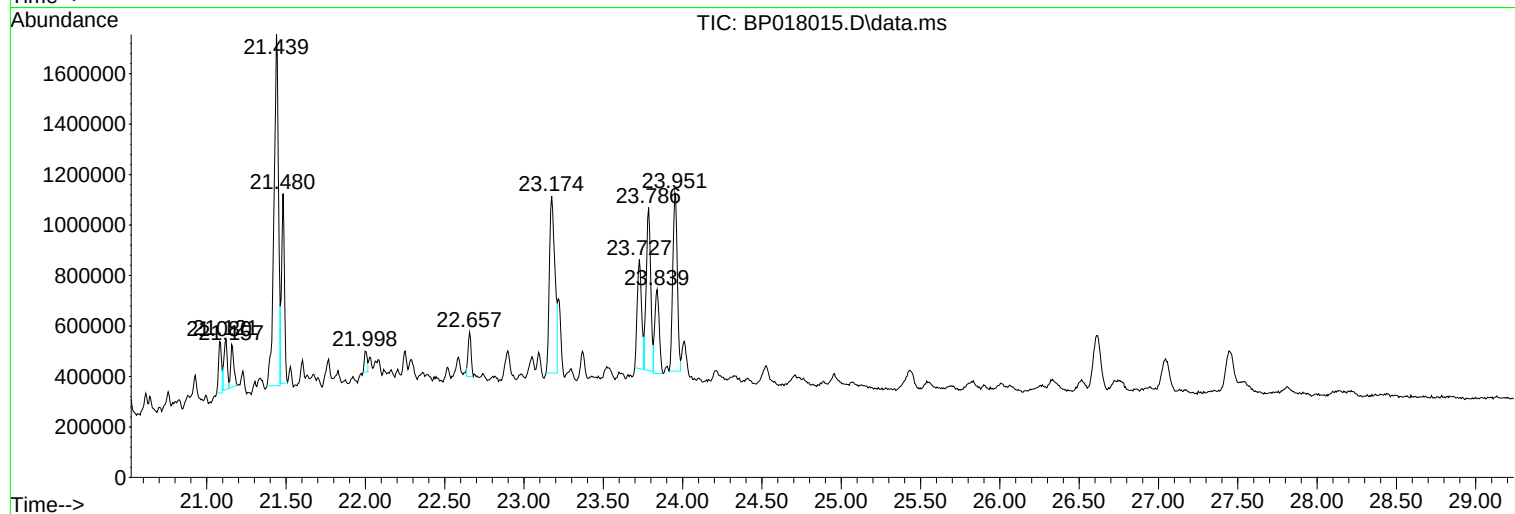
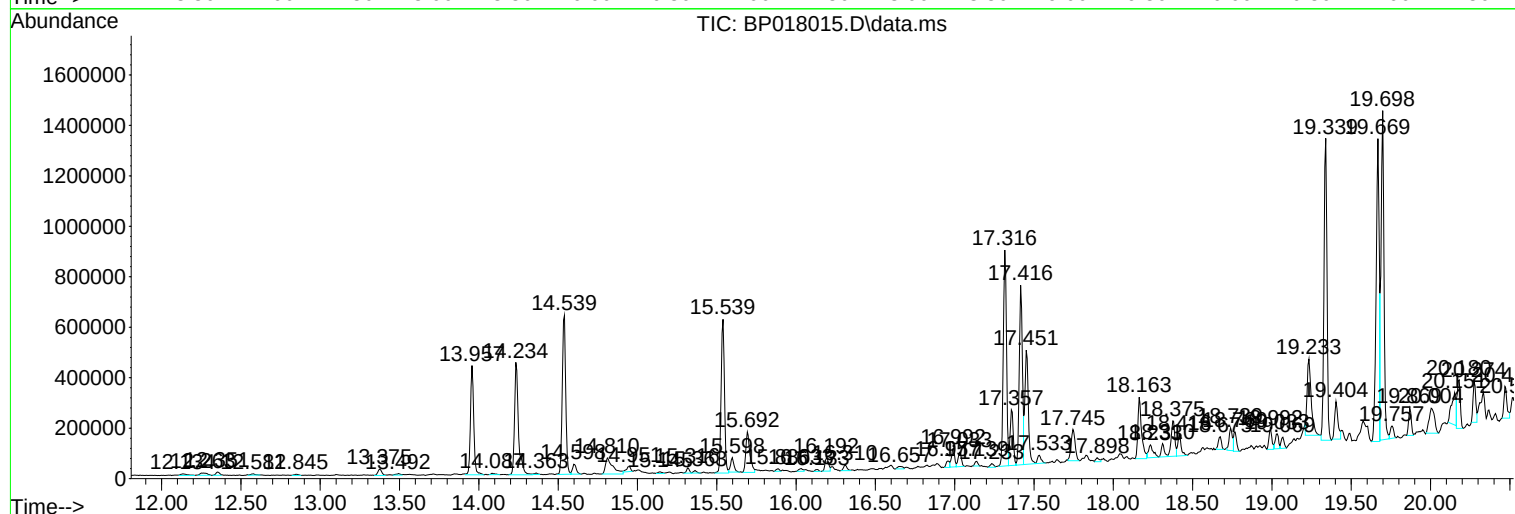
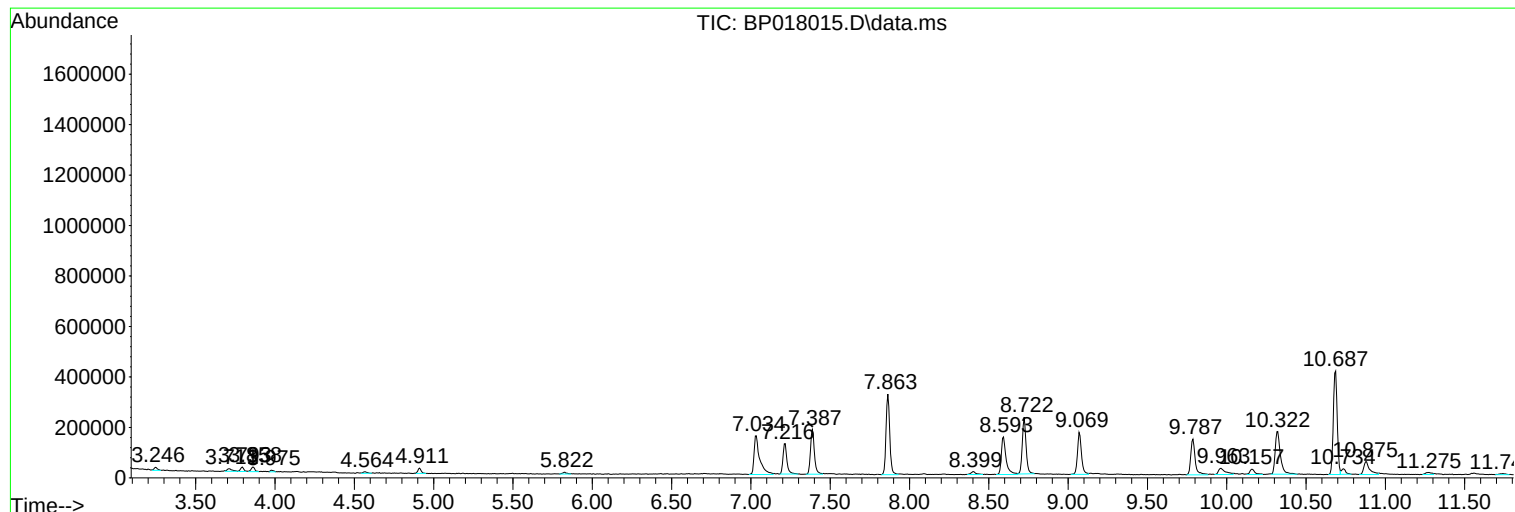
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110923.MA.M
Quant Title : SVOA CALIBRATION

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



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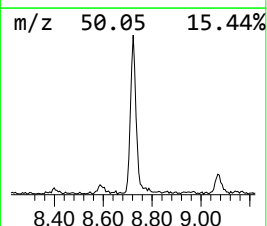
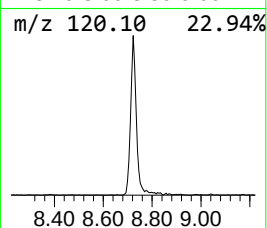
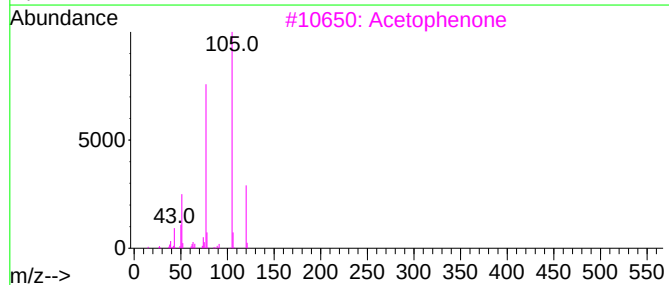
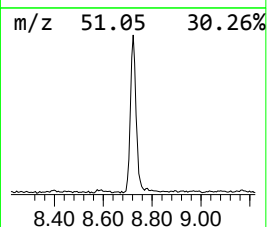
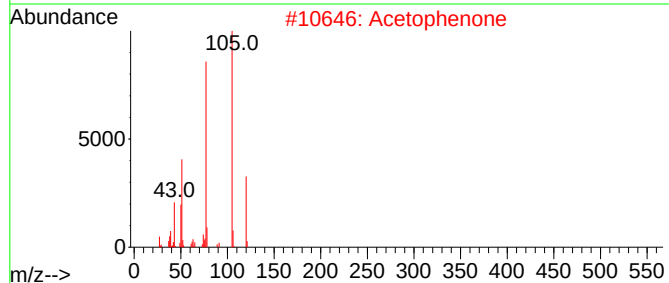
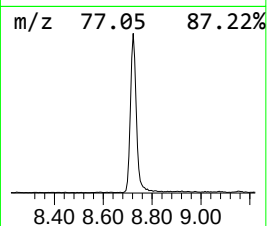
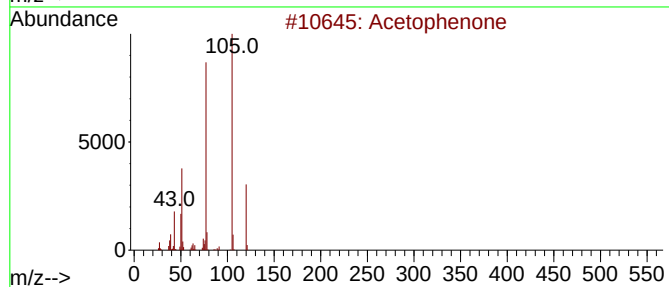
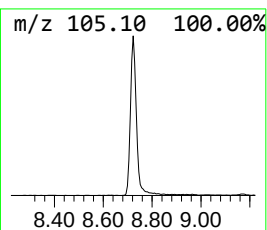
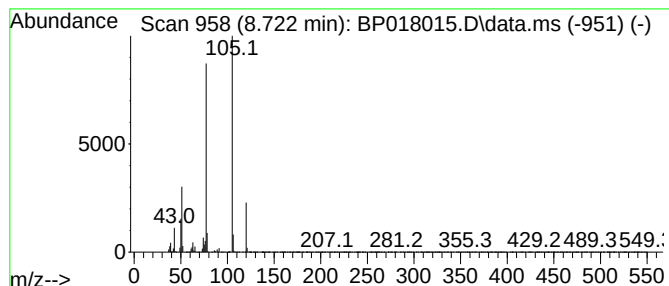
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110923.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Aceto phenone Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.722	14.39 ng/ul	372368	1,4-Dichlorobenzene-d4	7.863

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetophenone			120	C8H8O	000098-86-2	97
2	Acetophenone			120	C8H8O	000098-86-2	96
3	Acetophenone			120	C8H8O	000098-86-2	94
4	Acetophenone			120	C8H8O	000098-86-2	91
5	Acetophenone			120	C8H8O	000098-86-2	91



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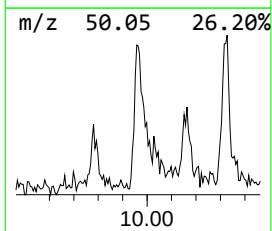
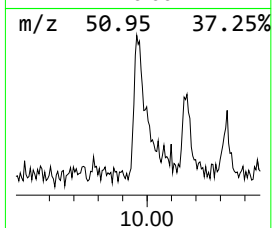
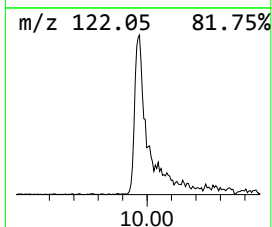
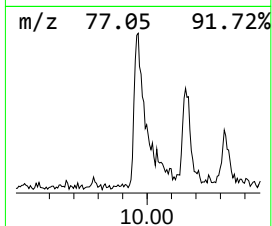
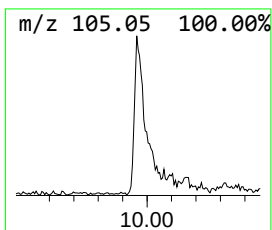
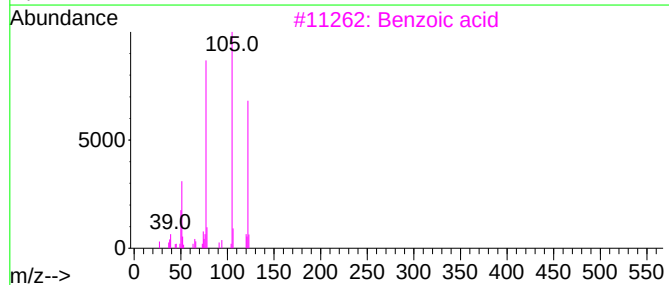
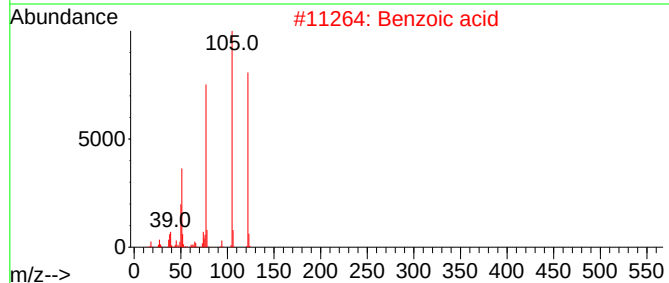
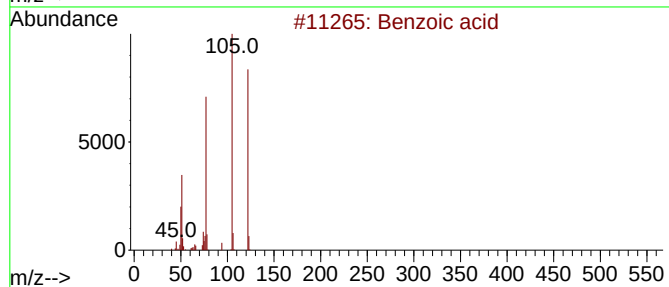
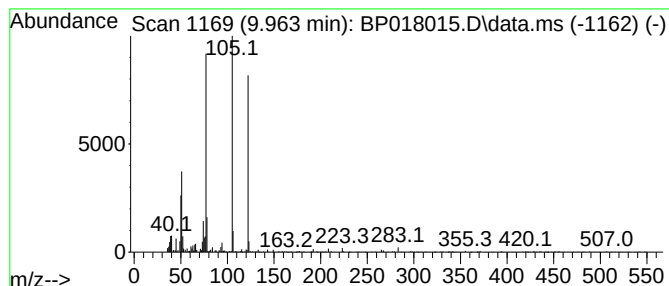
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110923.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.963	2.16 ng/ul	75857	Naphthalene-d8	10.687

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzoic acid	122	C7H6O2	000065-85-0	95
2			Benzoic acid	122	C7H6O2	000065-85-0	94
3			Benzoic acid	122	C7H6O2	000065-85-0	91
4			Benzoic acid	122	C7H6O2	000065-85-0	91
5			Cyclobutane-1,1-dicarboxamide, N...	382	C20H18N2O6	1000253-25-3	83



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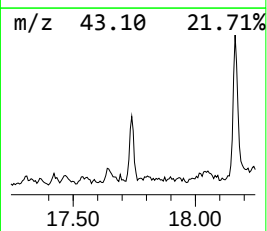
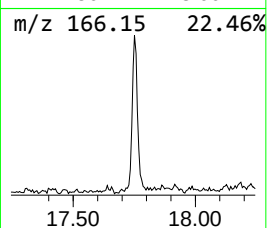
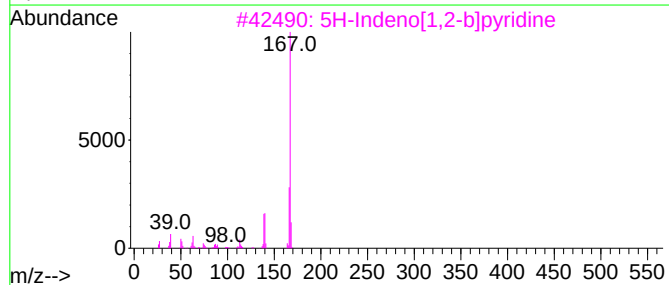
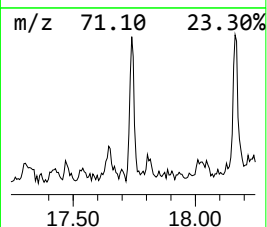
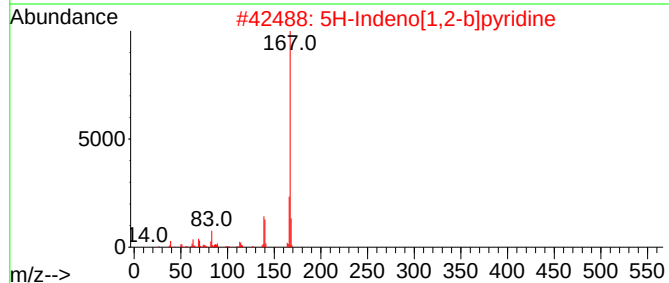
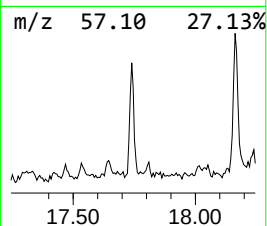
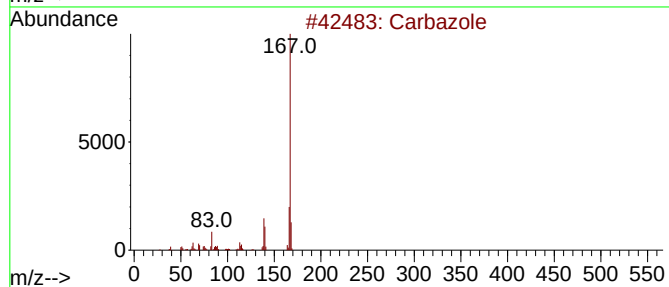
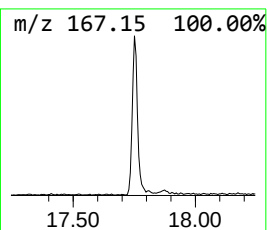
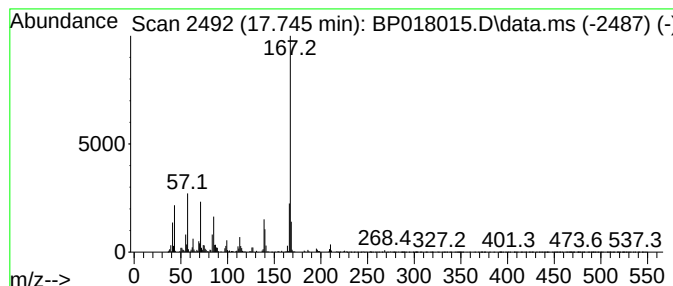
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110923.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Carba zole Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.745	3.61 ng/ul	217543	Phenanthrene-d10	17.316

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Carbazole	167	C12H9N	000086-74-8	93
2			5H-Indeno[1,2-b]pyridine	167	C12H9N	000244-99-5	76
3			5H-Indeno[1,2-b]pyridine	167	C12H9N	000244-99-5	64
4			Carbazole	167	C12H9N	000086-74-8	64
5			9H-Carbazole-9-methanol	197	C13H11NO	002409-36-1	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110923\
Data File : BP018015.D
Acq On : 10 Nov 2023 06:51
Operator : MA/JU
Sample : 05091-01
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCK10

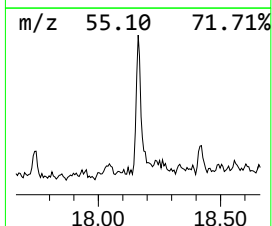
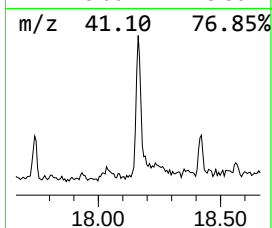
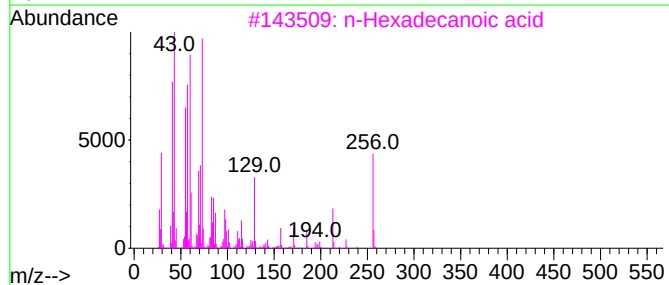
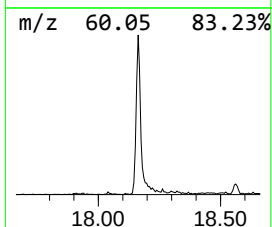
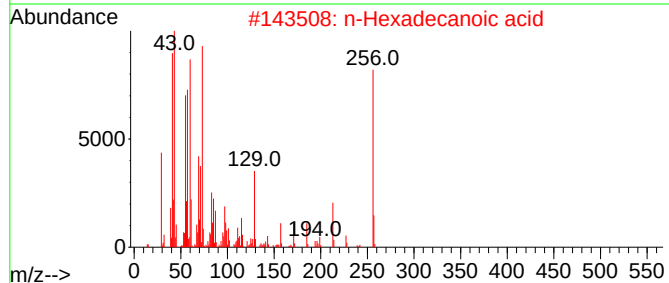
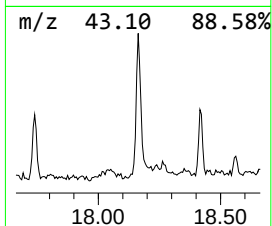
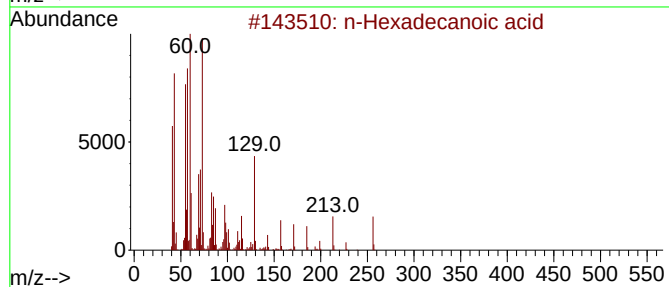
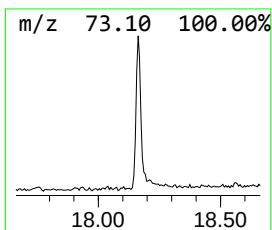
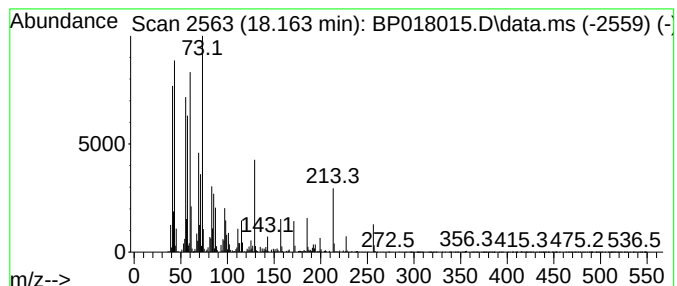
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.163	6.55 ng/ul	394575	Phenanthrene-d10	17.316

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	98
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
5			Octadecanoic acid	284	C18H36O2	000057-11-4	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110923\
Data File : BP018015.D
Acq On : 10 Nov 2023 06:51
Operator : MA/JU
Sample : 05091-01
Misc :
ALS Vial : 22 Sample Multiplier: 1

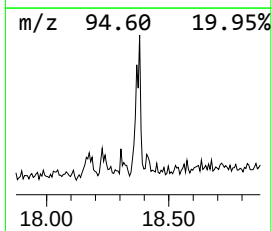
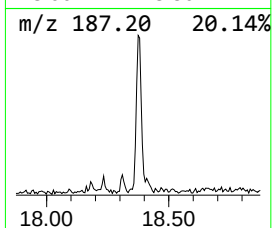
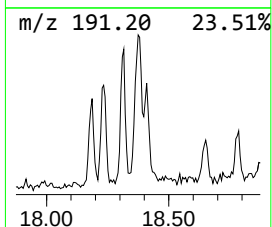
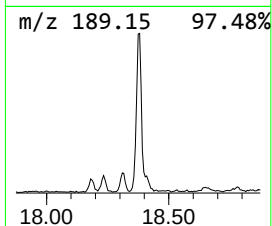
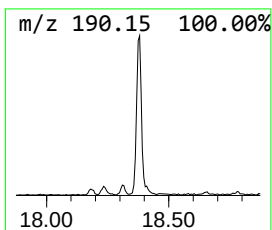
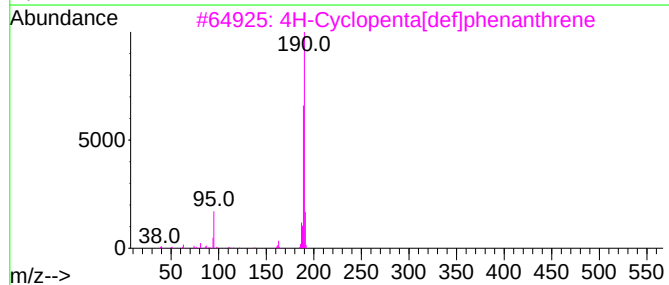
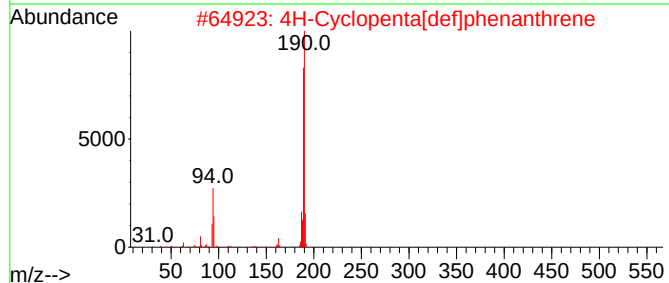
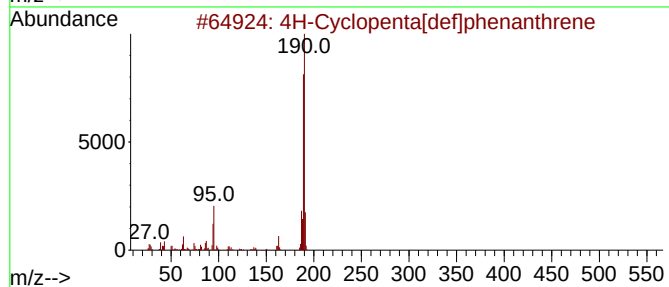
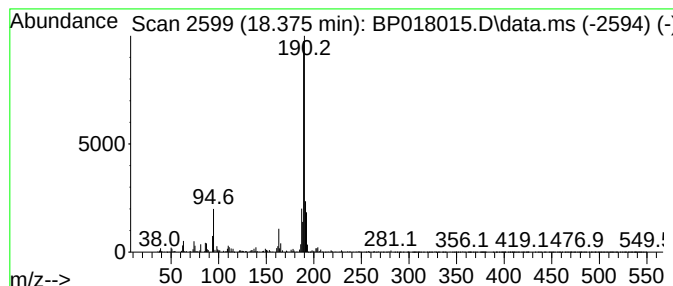
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 4H-Cyclopenta[def]phenanthrene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.375	3.79 ng/ul	228270	Phenanthrene-d10	17.316	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	95
2	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	87
3	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	72
4	Methyl diselenide	190	C2H6Se2	007101-31-7	49
5	2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	45



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110923\
Data File : BP018015.D
Acq On : 10 Nov 2023 06:51
Operator : MA/JU
Sample : 05091-01
Misc :
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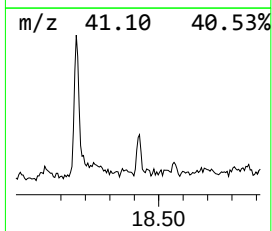
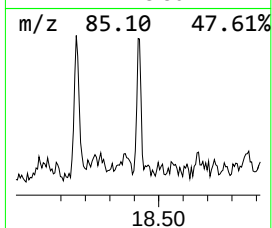
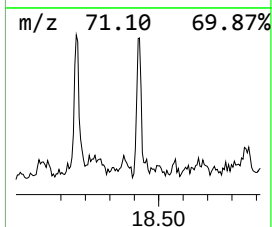
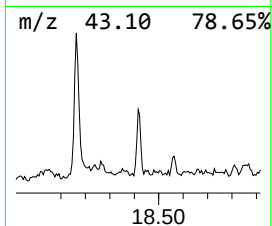
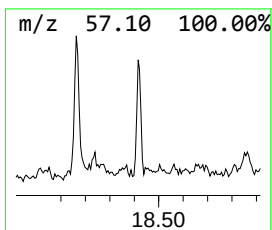
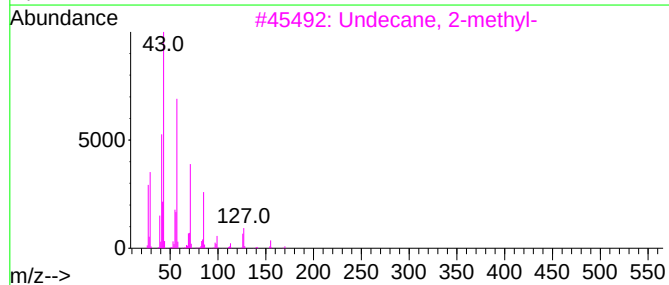
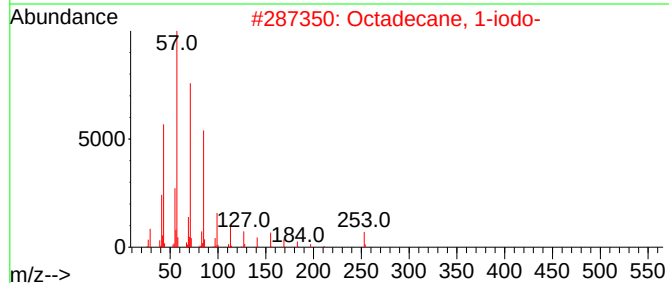
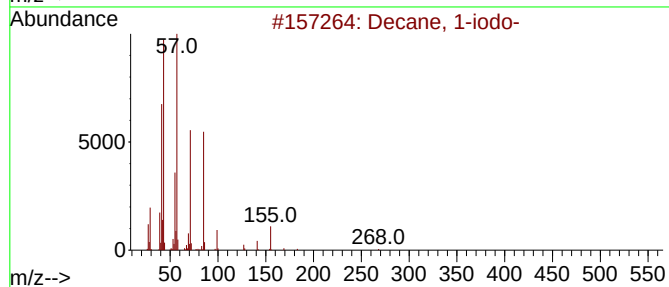
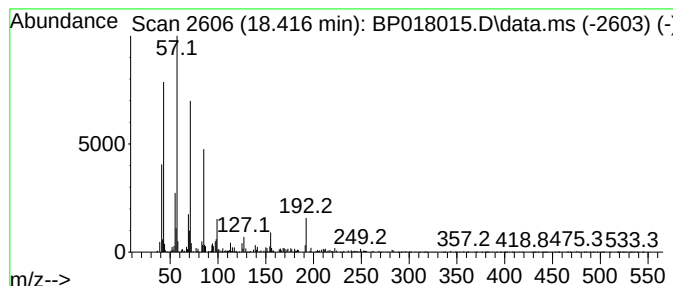
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Decane, 1-iodo- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD		R.T.	
18.416	2.06 ng/ul	124280	Phenanthrene-d10		17.316	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Decane, 1-iodo-		268	C10H21I	002050-77-3	87
2	Octadecane, 1-iodo-		380	C18H37I	000629-93-6	80
3	Undecane, 2-methyl-		170	C12H26	007045-71-8	80
4	Octadecane, 1-iodo-		380	C18H37I	000629-93-6	80
5	Dodecane, 1-iodo-		296	C12H25I	004292-19-7	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110923\
Data File : BP018015.D
Acq On : 10 Nov 2023 06:51
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Sample : 05091-01
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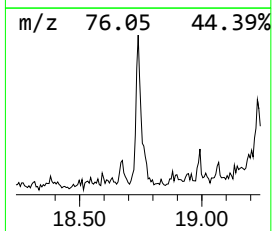
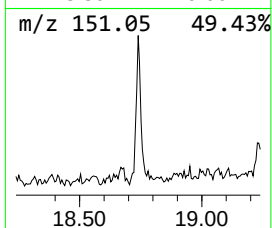
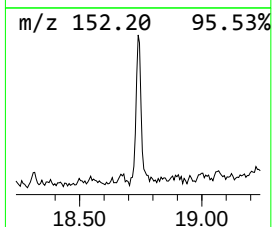
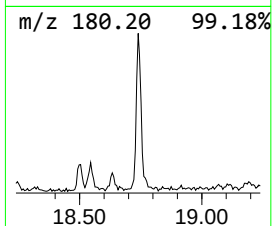
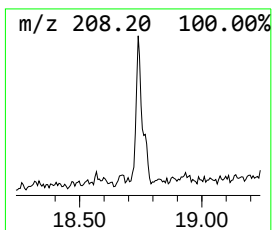
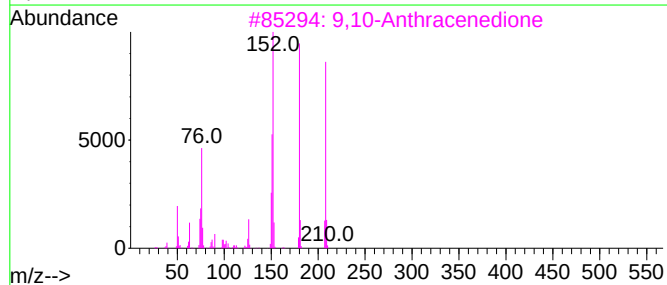
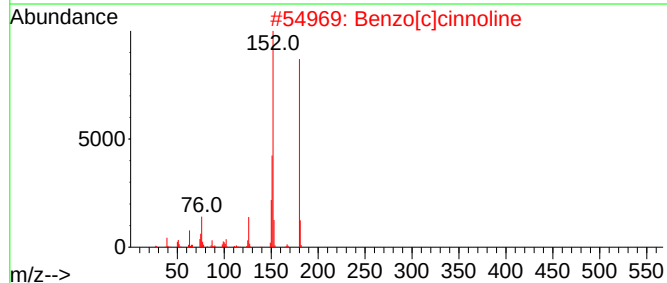
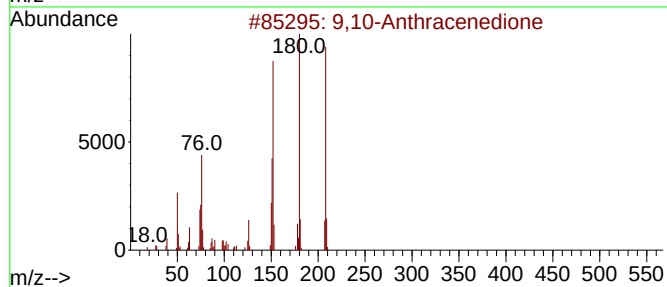
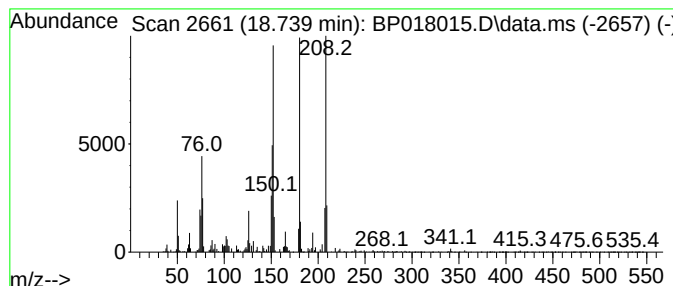
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 9,10-Anthracenedione Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.739	2.26 ng/ul	135939	Phenanthrene-d10	17.316

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9,10-Anthracenedione	208	C14H8O2	000084-65-1	95
2			Benzo[c]cinnoline	180	C12H8N2	000230-17-1	95
3			9,10-Anthracenedione	208	C14H8O2	000084-65-1	95
4			9,10-Anthracenedione	208	C14H8O2	000084-65-1	93
5			9H-Fluoren-9-one	180	C13H8O	000486-25-9	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110923\
Data File : BP018015.D
Acq On : 10 Nov 2023 06:51
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Sample : 05091-01
Misc :
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Instrument :
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ClientSampleId :
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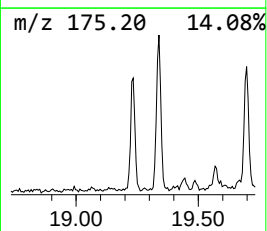
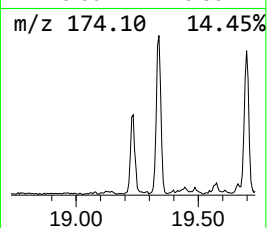
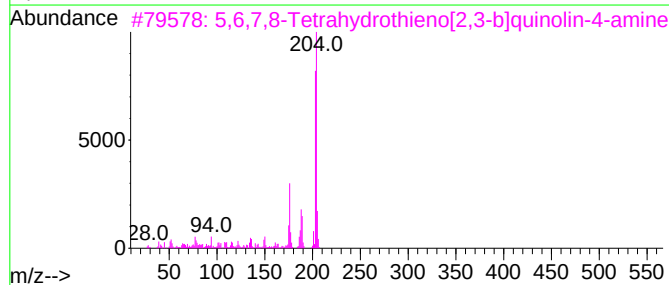
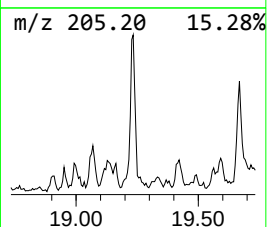
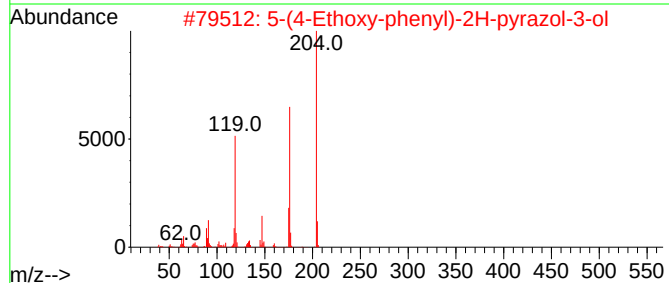
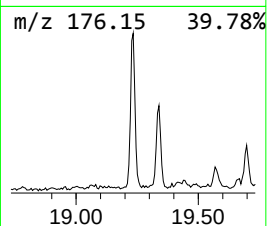
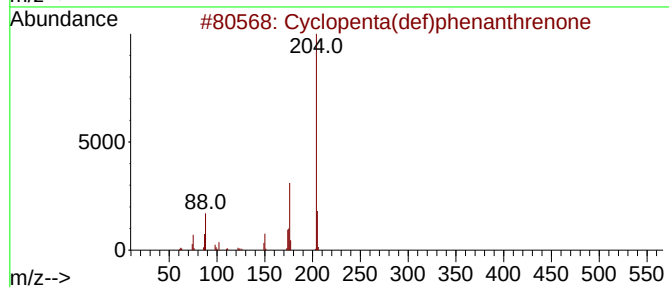
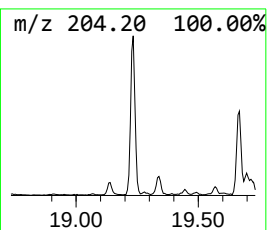
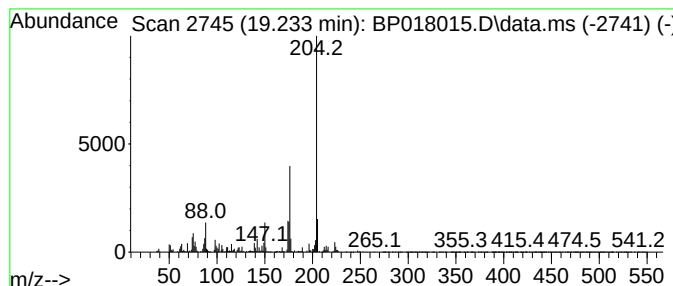
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Cyclopenta(def)phenanthrenone Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.233	9.05 ng/ul	545297	Phenanthrene-d10	17.316

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopenta(def)phenanthrenone	204	C15H8O	005737-13-3	93
2			5-(4-Ethoxy-phenyl)-2H-pyrazol-3-ol	204	C11H12N2O2	1000278-26-8	59
3			5,6,7,8-Tetrahydrothieno[2,3-b]q...	204	C11H12N2S	122914-50-5	50
4			[1,1'-Biphenyl]-2,2'-dicarbonitrile	204	C14H8N2	004341-02-0	50
5			Tricyclo[5.1.0.0(2,4)]octane, 5-...	247	C17H29N	1000063-59-3	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110923\
Data File : BP018015.D
Acq On : 10 Nov 2023 06:51
Operator : MA/JU
Sample : 05091-01
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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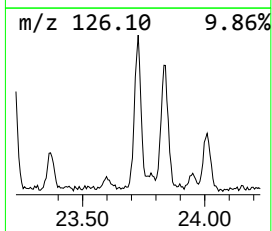
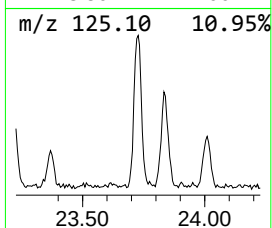
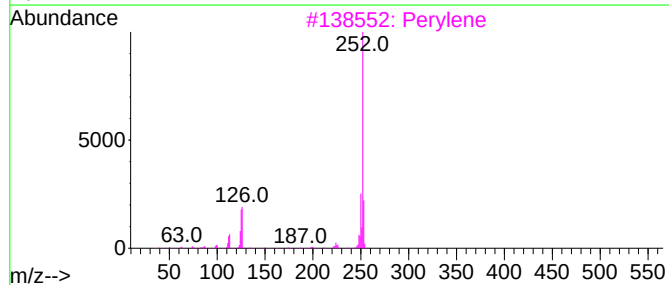
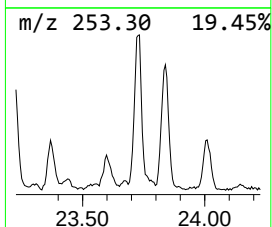
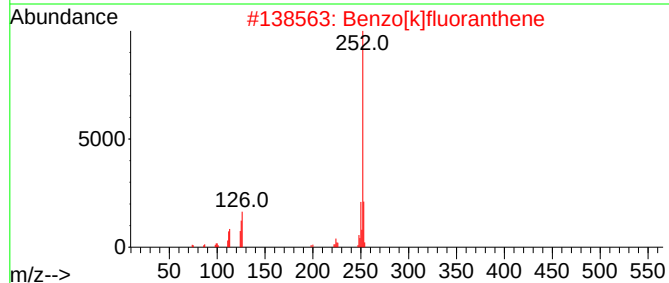
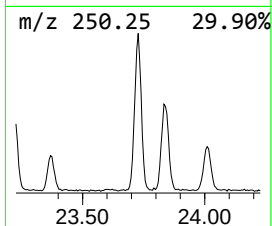
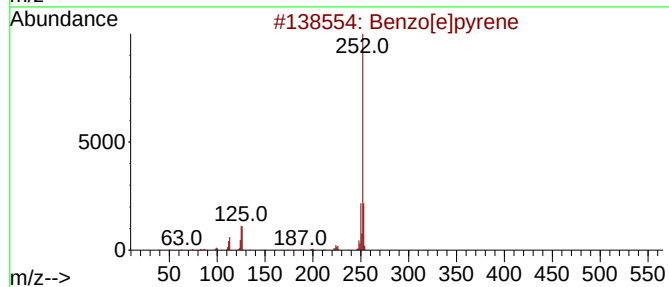
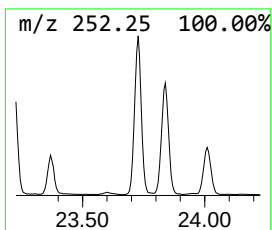
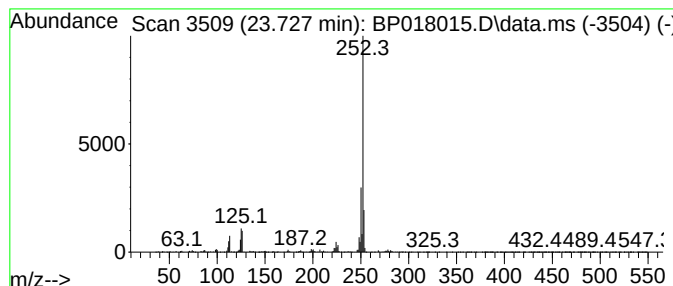
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Benzo[e]pyrene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.727	12.14 ng/ul	854837	Perylene-d12	23.951

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	97
3			Perylene	252	C20H12	000198-55-0	94
4			Benzo[j]fluoranthene	252	C20H12	000205-82-3	94
5			Benzo[b]fluoranthene	252	C20H12	000205-99-2	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110923\
Data File : BP018015.D
Acq On : 10 Nov 2023 06:51
Operator : MA/JU
Sample : 05091-01
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCK10

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110923.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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Benzoic acid	9.963	2.2	ng/ul	75857	2	10.687	701208	20.0
Carba zole	17.745	3.6	ng/ul	217543	4	17.316	1204710	20.0
n-Hexadecanoic ...	18.163	6.5	ng/ul	394575	4	17.316	1204710	20.0
4H-Cyclopenta[d]...	18.375	3.8	ng/ul	228270	4	17.316	1204710	20.0
Decane, 1-iodo-	18.416	2.1	ng/ul	124280	4	17.316	1204710	20.0
9,10-Anthracene...	18.739	2.3	ng/ul	135939	4	17.316	1204710	20.0
Cyclopenta(def)...	19.233	9.1	ng/ul	545297	4	17.316	1204710	20.0
Benzo[e]pyrene	23.727	12.1	ng/ul	854837	6	23.951	1408710	20.0