

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061721\
 Data File : BM030494.D
 Acq On : 17 Jun 2021 14:28
 Operator : CG/JU
 Sample : M2596-16DL 2X
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BG5R5DL

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-BM060621.M
 Title : SVOA CALIBRATION

Signal : TIC: BM030494.D

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.752	107	113	115	rBV	27481	48189	0.99%	0.087%
2	3.781	115	118	123	rVV2	36189	74961	1.54%	0.135%
3	3.840	123	128	137	rVB	151511	310948	6.40%	0.558%
4	4.057	160	165	176	rVB2	29732	76629	1.58%	0.138%
5	4.604	253	258	266	rVB	54775	107568	2.21%	0.193%
6	4.963	314	319	329	rVB	25221	51032	1.05%	0.092%
7	5.063	331	336	345	rBV2	23384	53045	1.09%	0.095%
8	5.493	402	409	419	rBV	67398	166690	3.43%	0.299%
9	5.857	464	471	477	rBV2	36311	65978	1.36%	0.118%
10	6.998	657	665	675	rBV	179751	325396	6.70%	0.584%
11	7.163	685	693	703	rBV	156214	270625	5.57%	0.486%
12	7.363	720	727	736	rBV	201046	345079	7.10%	0.620%
13	7.828	798	806	814	rVV	861867	1399705	28.80%	2.513%
14	8.528	916	925	934	rBV	109270	200663	4.13%	0.360%
15	8.981	996	1002	1011	rVV	181274	323065	6.65%	0.580%
16	9.698	1116	1124	1130	rBV	146821	252313	5.19%	0.453%
17	10.028	1175	1180	1187	rVB4	29509	52145	1.07%	0.094%
18	10.239	1207	1216	1226	rBV	201718	386201	7.95%	0.693%
19	10.410	1238	1245	1261	rBV	33717	95485	1.96%	0.171%
20	10.610	1271	1279	1285	rBV	1072964	1819472	37.44%	3.267%
21	10.657	1285	1287	1294	rVV	66481	96758	1.99%	0.174%
22	10.751	1296	1303	1318	rBV	101496	210505	4.33%	0.378%
23	12.269	1555	1561	1568	rVB2	51822	82703	1.70%	0.148%
24	12.486	1594	1598	1605	rVB	48009	77793	1.60%	0.140%
25	13.274	1726	1732	1738	rBV2	28221	48282	0.99%	0.087%
26	13.345	1738	1744	1752	rVB2	34975	60362	1.24%	0.108%
27	13.769	1811	1816	1821	rBV	45250	77105	1.59%	0.138%
28	13.857	1821	1831	1838	rVV	407250	647327	13.32%	1.162%
29	14.004	1848	1856	1860	rBV4	25536	43047	0.89%	0.077%
30	14.139	1873	1879	1893	rVB	491361	800948	16.48%	1.438%
31	14.445	1922	1931	1938	rVV	1490657	2252527	46.35%	4.044%
32	14.510	1938	1942	1949	rVV	150571	231782	4.77%	0.416%
33	14.657	1961	1967	1976	rVV3	76781	169075	3.48%	0.304%
34	14.751	1978	1983	1989	rVV3	20192	42400	0.87%	0.076%
35	14.845	1995	1999	2006	rVV2	86444	157180	3.23%	0.282%
36	15.439	2091	2100	2105	rVV	646774	942736	19.40%	1.693%

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Integrator: RTE

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Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM060621.M
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37	15.492	2105	2109	2116	rVV	165314	247815	5.10%	0.445%
38	15.557	2116	2120	2125	rVV	89736	148056	3.05%	0.266%
39	15.616	2128	2130	2133	rVV2	33492	40344	0.83%	0.072%
40	15.657	2133	2137	2142	rVV4	31503	54492	1.12%	0.098%
41	15.786	2155	2159	2166	rVB	44297	69461	1.43%	0.125%
42	15.857	2166	2171	2175	rBV	153357	199653	4.11%	0.358%
43	15.933	2180	2184	2192	rVB2	39349	85393	1.76%	0.153%
44	16.163	2218	2223	2229	rVB5	33853	67236	1.38%	0.121%
45	16.492	2268	2279	2284	rBV4	45696	134299	2.76%	0.241%
46	16.545	2284	2288	2293	rVB3	32078	49308	1.01%	0.089%
47	16.645	2301	2305	2310	rVB5	25011	41327	0.85%	0.074%
48	16.839	2333	2338	2345	rVB	53619	87110	1.79%	0.156%
49	16.945	2353	2356	2362	rVV5	66607	129216	2.66%	0.232%
50	17.004	2362	2366	2374	rVB	87985	144706	2.98%	0.260%
51	17.186	2391	2397	2401	rBV	1734013	2541797	52.30%	4.563%
52	17.227	2401	2404	2409	rVV	1816508	2402026	49.42%	4.312%
53	17.286	2409	2414	2417	rVV	641842	1012848	20.84%	1.818%
54	17.315	2417	2419	2425	rVV	368112	467807	9.63%	0.840%
55	17.380	2425	2430	2434	rVV2	40821	68703	1.41%	0.123%
56	17.586	2458	2465	2470	rBV2	112191	198332	4.08%	0.356%
57	17.704	2482	2485	2491	rVB	39209	57647	1.19%	0.103%
58	17.762	2492	2495	2502	rBV2	37918	59035	1.21%	0.106%
59	17.909	2516	2520	2526	rVB2	20664	38668	0.80%	0.069%
60	18.062	2539	2546	2550	rBV	257638	511659	10.53%	0.919%
61	18.104	2550	2553	2556	rVV	323846	472488	9.72%	0.848%
62	18.151	2556	2561	2565	rVV	3544639	4567570	93.98%	8.200%
63	18.186	2565	2567	2572	rVV2	162267	209828	4.32%	0.377%
64	18.257	2573	2579	2582	rVV2	341083	637173	13.11%	1.144%
65	18.292	2582	2585	2591	rVB	181125	260996	5.37%	0.469%
66	18.368	2595	2598	2601	rVB2	41200	51868	1.07%	0.093%
67	18.556	2626	2630	2634	rVB	158509	208232	4.28%	0.374%
68	18.598	2634	2637	2645	rVB2	111508	178531	3.67%	0.321%
69	18.680	2649	2651	2658	rVB	31044	40103	0.83%	0.072%
70	18.804	2669	2672	2675	rVB	61448	66476	1.37%	0.119%
71	18.862	2676	2682	2686	rBV	711563	1132549	23.30%	2.033%
72	18.974	2697	2701	2705	rBV	139106	233980	4.81%	0.420%
73	19.115	2721	2725	2735	rBV3	105717	227103	4.67%	0.408%
74	19.239	2741	2746	2753	rVB	2646612	3349010	68.91%	6.013%
75	19.304	2754	2757	2759	rBV2	46495	48407	1.00%	0.087%
76	19.339	2759	2763	2768	rBV5	77196	141504	2.91%	0.254%

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 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	19.574	2797	2803	2805	rBV3	1088152	1713436	35.26%	3.076%
78	19.603	2805	2808	2816	rVB2	2328367	3445280	70.89%	6.185%
79	19.674	2816	2820	2825	rBV3	124647	221610	4.56%	0.398%
80	19.921	2858	2862	2874	rVB4	152483	446929	9.20%	0.802%
81	20.062	2881	2886	2887	rBV4	118169	182772	3.76%	0.328%
82	20.092	2888	2891	2899	rVB	309035	481766	9.91%	0.865%
83	20.192	2905	2908	2912	rBV2	182668	225726	4.64%	0.405%
84	20.251	2913	2918	2922	rBV2	205817	333630	6.86%	0.599%
85	20.292	2922	2925	2928	rBV	98426	133201	2.74%	0.239%
86	20.386	2936	2941	2945	rBV2	137774	283637	5.84%	0.509%
87	20.433	2947	2949	2953	rVB	132500	160035	3.29%	0.287%
88	20.515	2960	2963	2965	rBV	134489	155991	3.21%	0.280%
89	20.545	2965	2968	2974	rVB	636603	844423	17.38%	1.516%
90	20.615	2977	2980	2984	rVB2	165160	208464	4.29%	0.374%
91	20.839	3015	3018	3024	rVB	181713	251432	5.17%	0.451%
92	21.133	3064	3068	3076	rVB2	400143	545542	11.23%	0.979%
93	21.268	3088	3091	3094	rVB	378375	386246	7.95%	0.693%
94	21.350	3099	3105	3109	rVV2	2605768	4859957	100.00%	8.725%
95	21.392	3109	3112	3118	rVB	1168949	1504282	30.95%	2.701%
96	21.503	3128	3131	3137	rVB2	217674	302237	6.22%	0.543%
97	21.674	3156	3160	3165	rVB2	808280	1017033	20.93%	1.826%
98	22.203	3248	3250	3257	rVB2	273013	355424	7.31%	0.638%
99	22.944	3371	3376	3381	rBV	862018	1861002	38.29%	3.341%
100	23.633	3487	3493	3500	rBV2	1487987	2731548	56.21%	4.904%

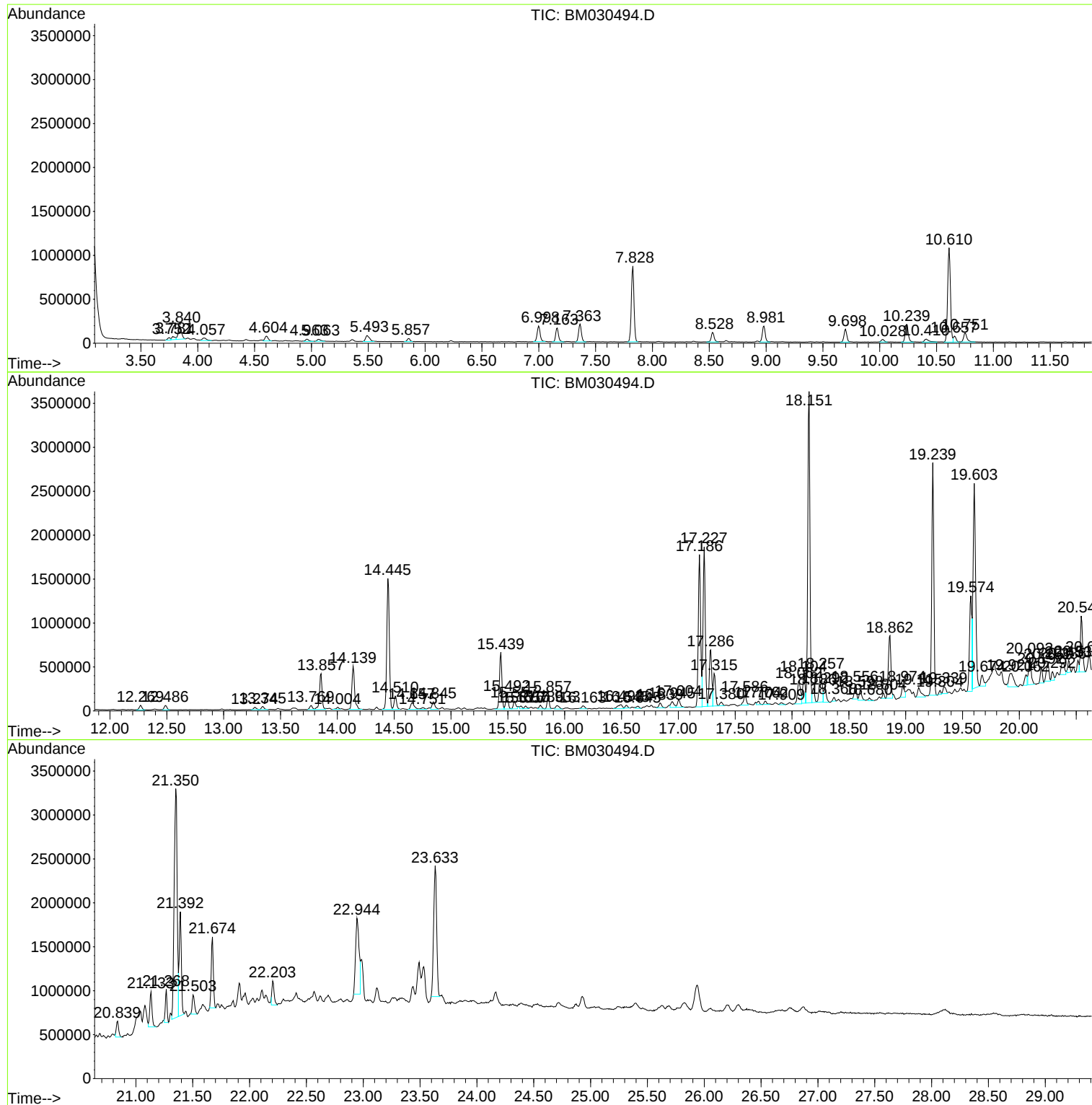
Sum of corrected areas: 55700078

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM060621.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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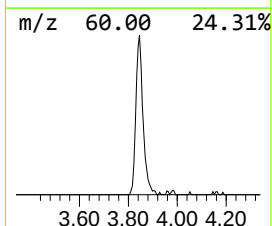
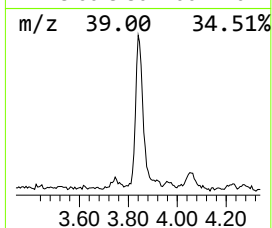
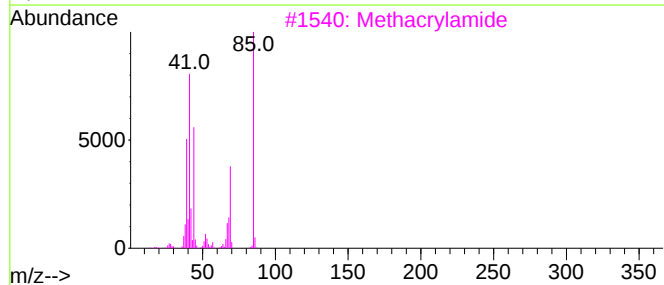
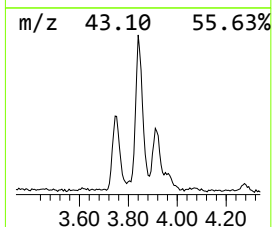
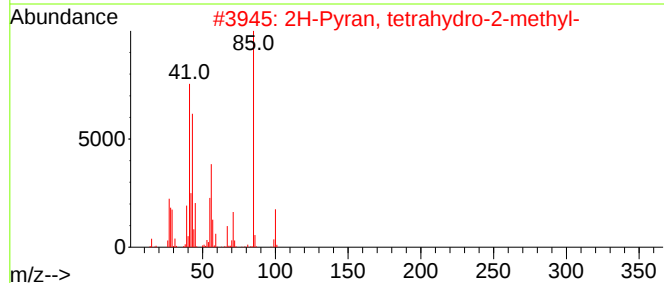
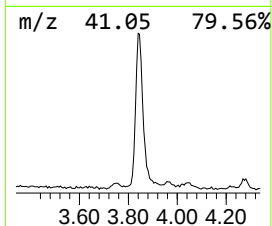
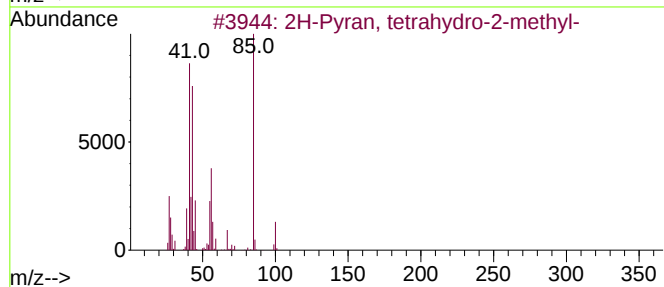
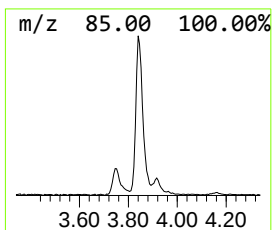
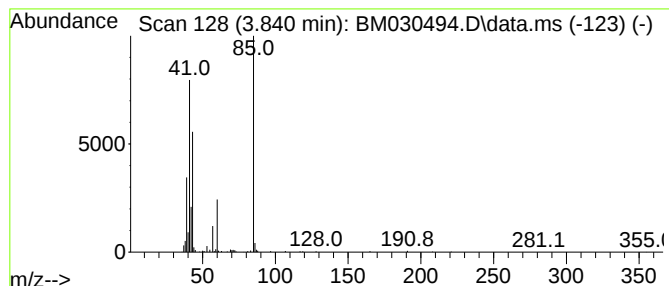
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM060621.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 2H-Pyran, tetrahydro-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.840	4.44 ng/ul	310948	1,4-Dichlorobenzene-d4	7.828

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2H-Pyran, tetrahydro-2-methyl-			100	C6H12O	010141-72-7	72
2	2H-Pyran, tetrahydro-2-methyl-			100	C6H12O	010141-72-7	38
3	Methacrylamide			85	C4H7NO	000079-39-0	36
4	Methacrylamide			85	C4H7NO	000079-39-0	36
5	Thiazole			85	C3H3NS	000288-47-1	9



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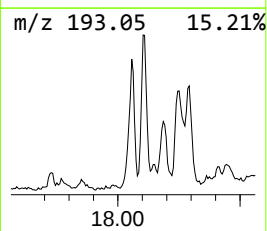
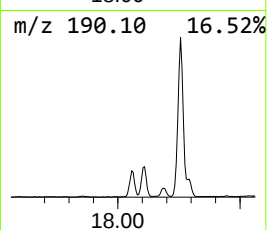
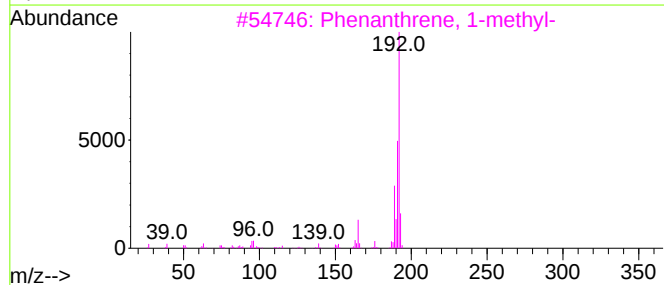
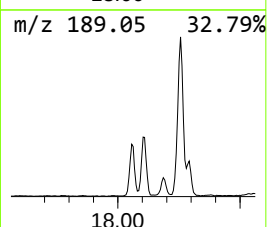
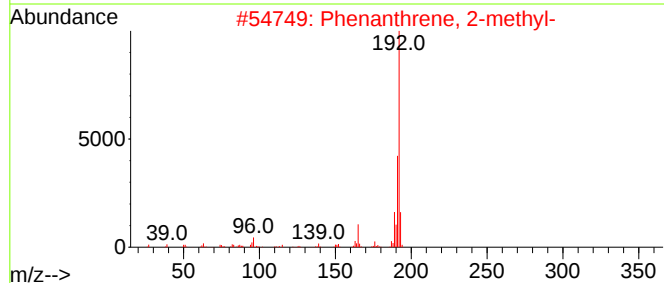
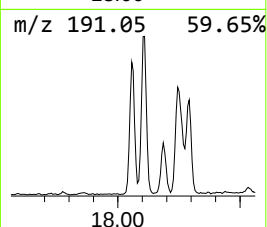
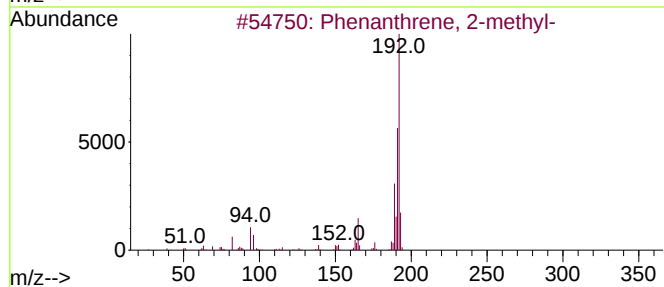
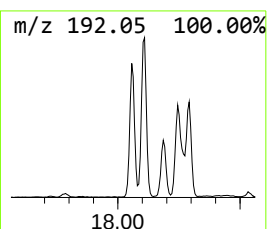
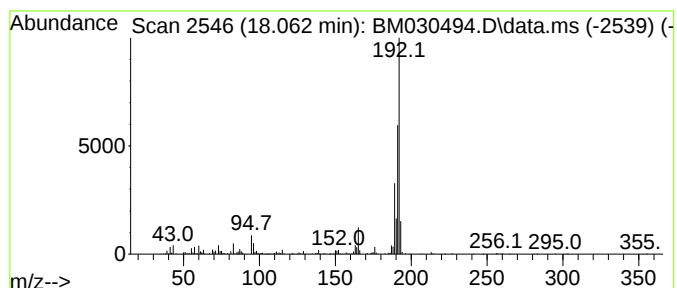
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Phenanthrene, 2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.060	4.03 ng/ul	511659	Phenanthrene-d10	17.186

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	97
2	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
3	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95
4	Anthracene, 2-methyl-	192	C15H12	000613-12-7	95
5	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	94



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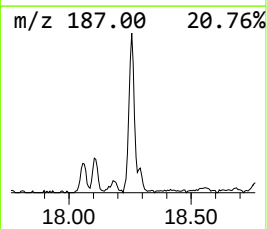
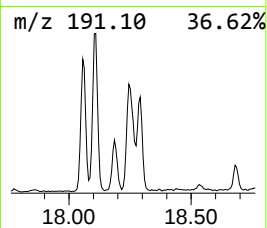
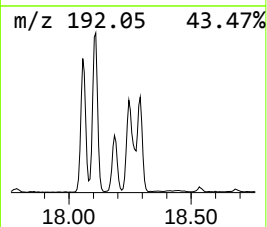
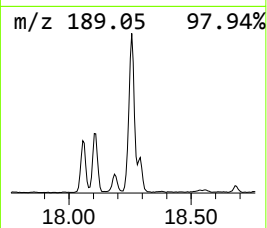
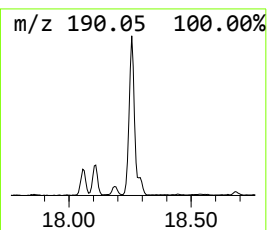
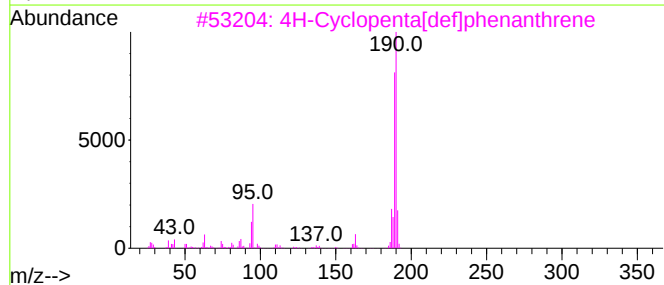
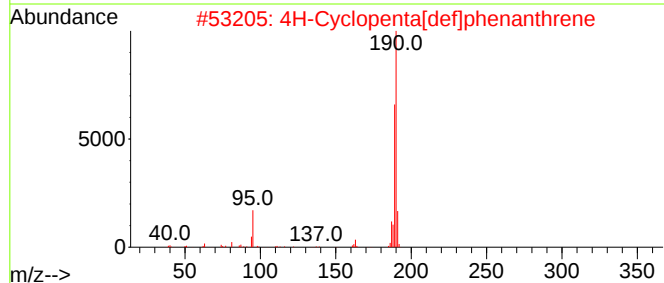
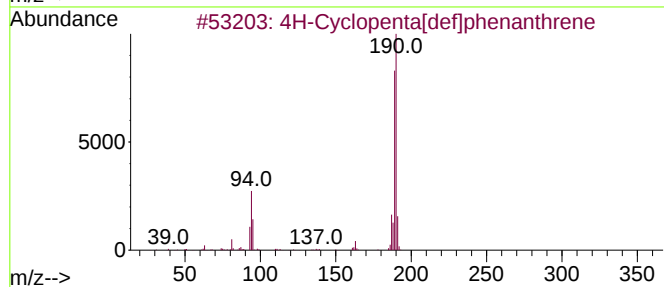
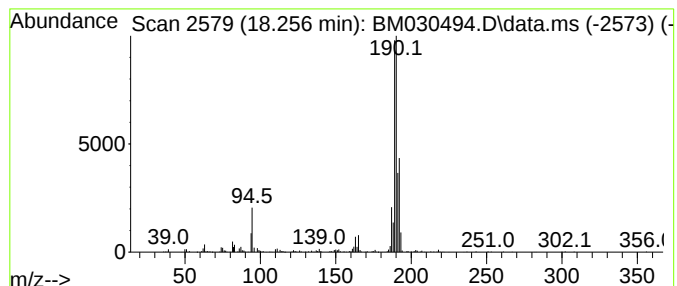
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TIC Integration Parameters: LSCINT.P

Peak Number 4 4H-Cyclopenta[def]phenanthrene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.260	5.01 ng/ul	637173	Phenanthrene-d10	17.186

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	94
2			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	70
3			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	52
4			2,3,5,6-Tetramethylterephthalald...	190	C12H14O2	007072-01-7	47
5			6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	42



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061721\
Data File : BM030494.D
Acq On : 17 Jun 2021 14:28
Operator : CG/JU
Sample : M2596-16DL 2X
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BG5R5DL

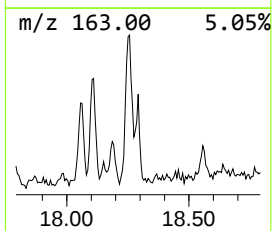
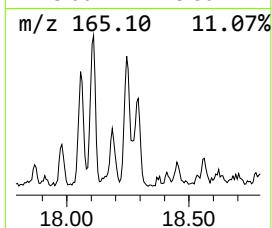
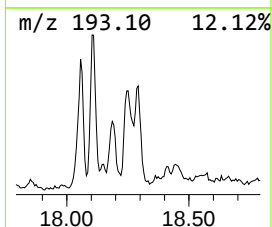
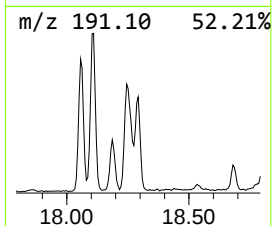
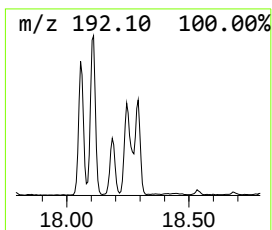
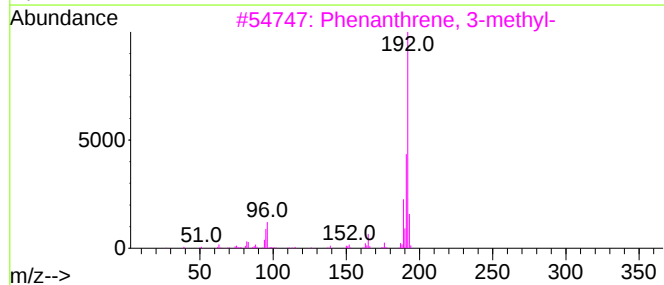
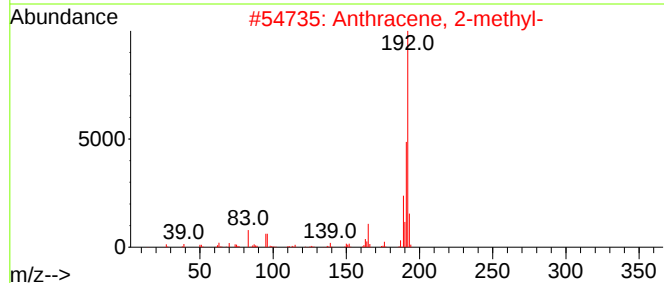
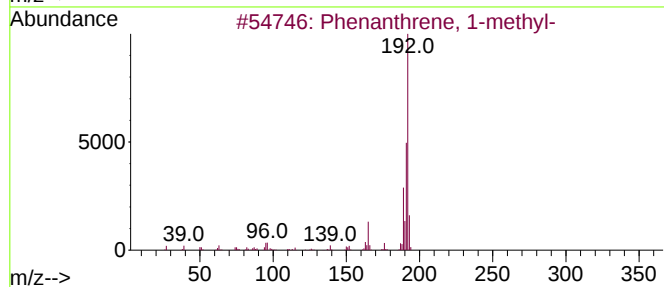
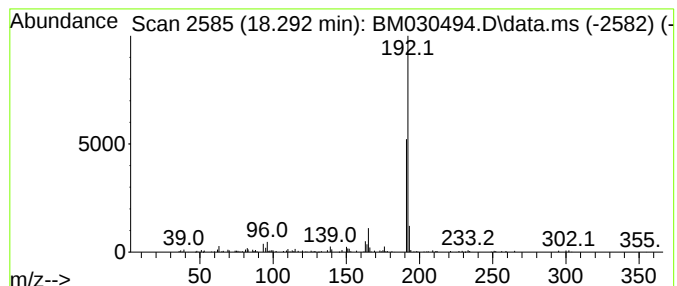
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Anthracene, 2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.290	2.05 ng/ul	260996	Phenanthrene-d10	17.186

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	94
2			Anthracene, 2-methyl-	192	C15H12	000613-12-7	94
3			Phenanthrene, 3-methyl-	192	C15H12	000832-71-3	91
4			Anthracene, 1-methyl-	192	C15H12	000610-48-0	91
5			Anthracene, 1-methyl-	192	C15H12	000610-48-0	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061721\
Data File : BM030494.D
Acq On : 17 Jun 2021 14:28
Operator : CG/JU
Sample : M2596-16DL 2X
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BG5R5DL

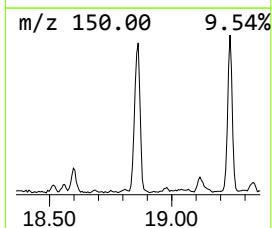
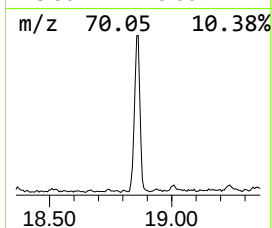
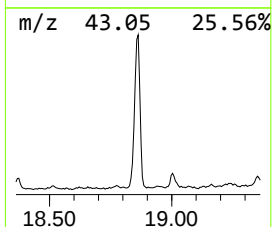
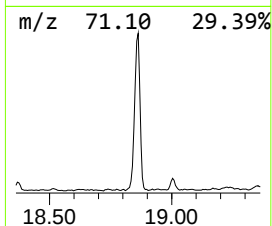
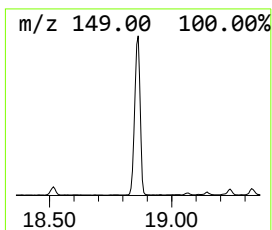
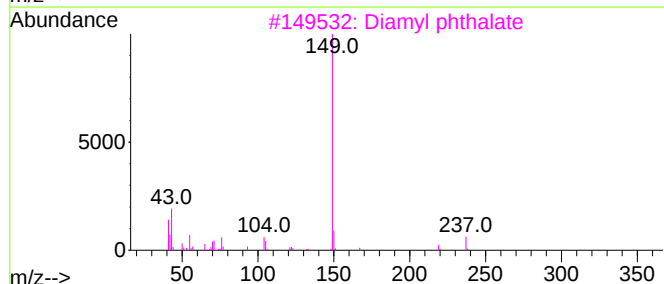
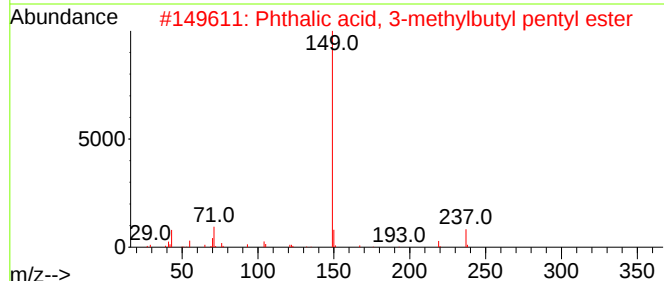
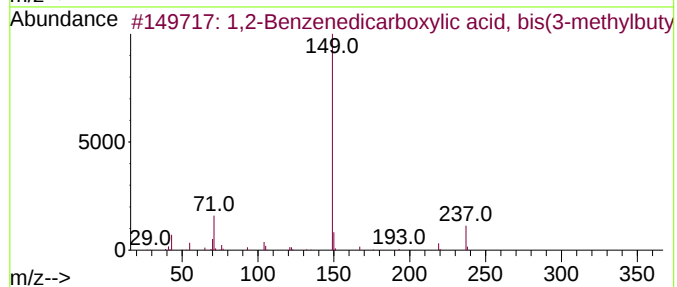
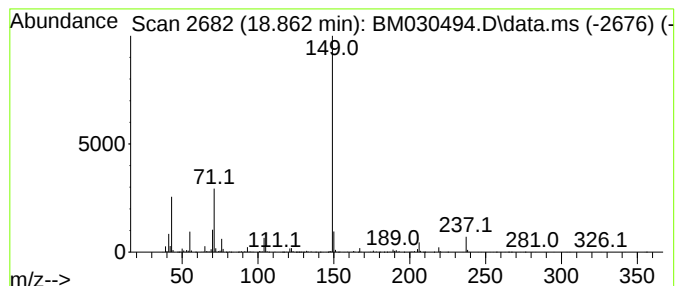
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 1,2-Benzenedicarboxylic aci... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.860	8.91 ng/ul	1132550	Phenanthrene-d10	17.186

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,2-Benzenedicarboxylic acid, bi...	306	C18H26O4	000605-50-5	90
2			Phthalic acid, 3-methylbutyl pen...	306	C18H26O4	1000356-80-5	83
3			Diamyl phthalate	306	C18H26O4	000131-18-0	72
4			Diamyl phthalate	306	C18H26O4	000131-18-0	72
5			Phthalic acid, 2-methylbutyl pen...	306	C18H26O4	1000322-81-2	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061721\
Data File : BM030494.D
Acq On : 17 Jun 2021 14:28
Operator : CG/JU
Sample : M2596-16DL 2X
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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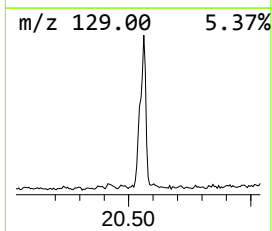
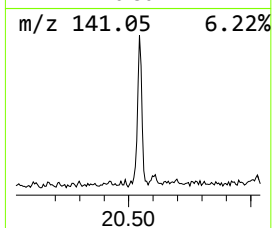
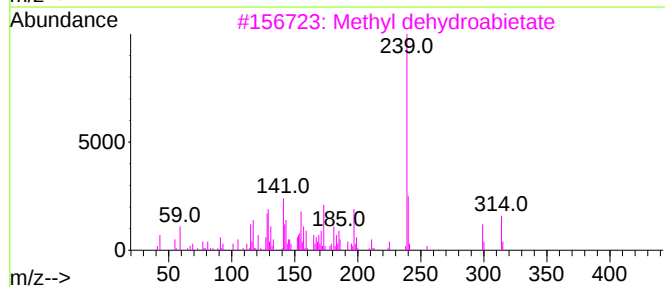
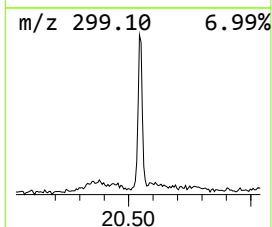
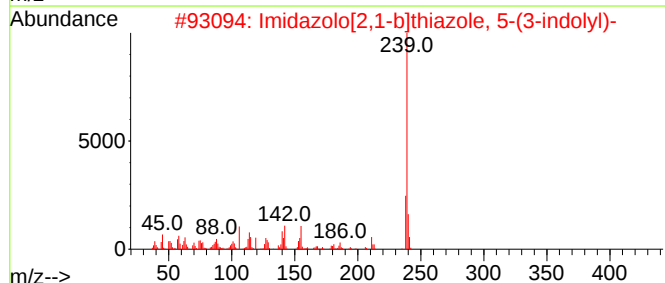
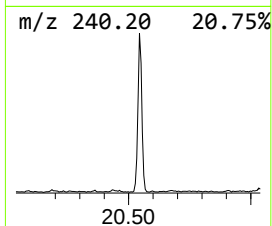
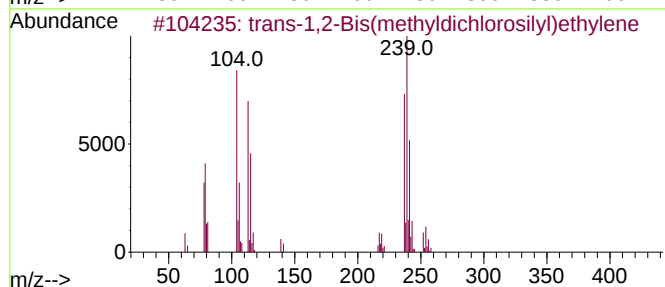
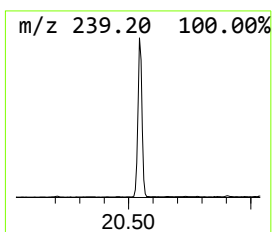
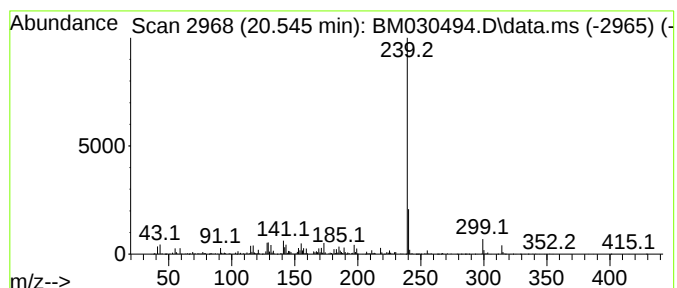
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 trans-1,2-Bis(methyldichlor... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.540	3.48 ng/ul	844423	Chrysene-d12	21.356

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			trans-1,2-Bis(methyldichlorosilyl)ethylen...	252	C4H8Cl4Si2	065899-10-7	60
2			Imidazolo[2,1-b]thiazole, 5-(3-indolyl)-	239	C13H9N3S	327097-37-0	59
3			Methyl dehydroabietate	314	C21H30O2	001235-74-1	56
4			Methyl dehydroabietate	314	C21H30O2	001235-74-1	55
5			4-(4-Oxo-1,2,3,4,6,7,12,12b-octa...	326	C19H22N2O3	1000137-91-1	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061721\
Data File : BM030494.D
Acq On : 17 Jun 2021 14:28
Operator : CG/JU
Sample : M2596-16DL 2X
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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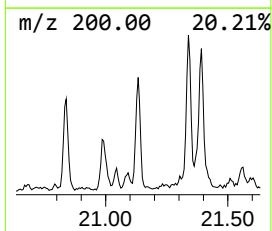
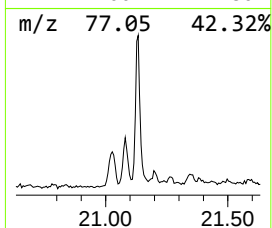
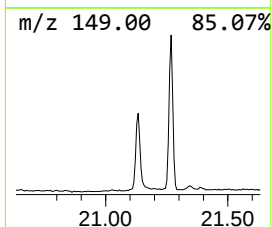
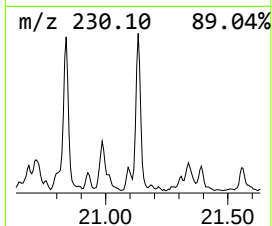
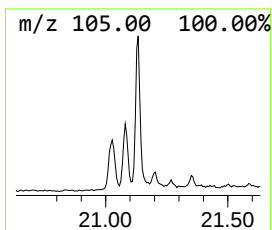
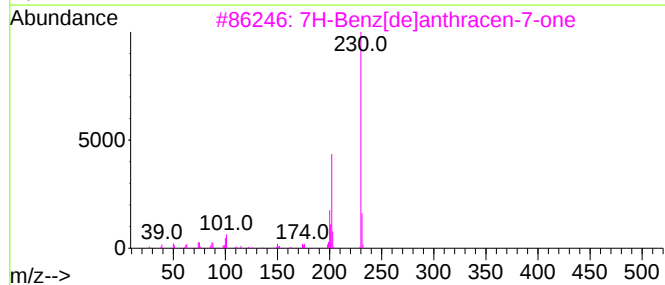
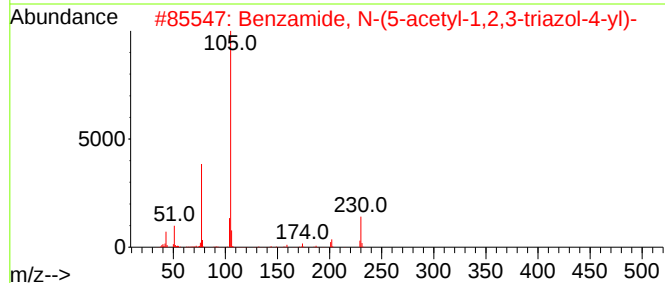
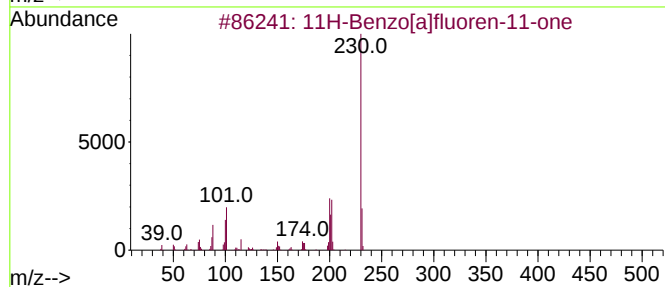
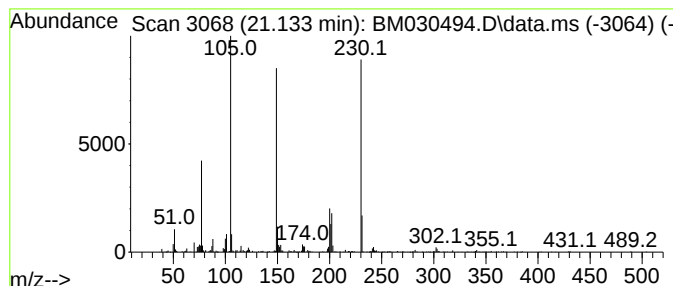
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 11H-Benzo[a]fluoren-11-one Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.130	2.25 ng/ul	545542	Chrysene-d12	21.356

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			11H-Benzo[a]fluoren-11-one	230	C17H10O	000479-79-8	87
2			Benzamide, N-(5-acetyl-1,2,3-tri...	230	C11H10N4O2	129959-33-7	53
3			7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	50
4			Diethylene glycol dibenzoate	314	C18H18O5	000120-55-8	43
5			7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061721\
Data File : BM030494.D
Acq On : 17 Jun 2021 14:28
Operator : CG/JU
Sample : M2596-16DL 2X
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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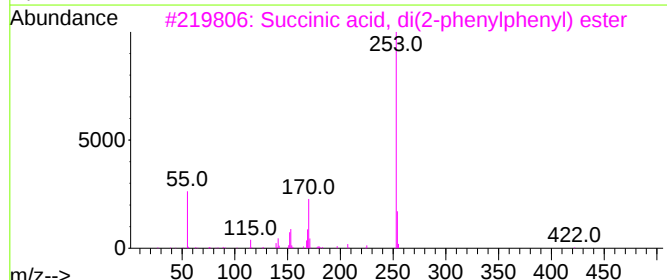
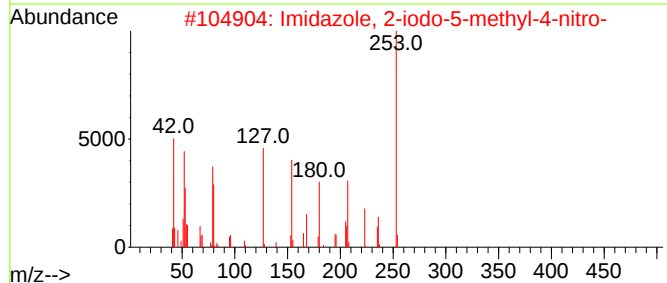
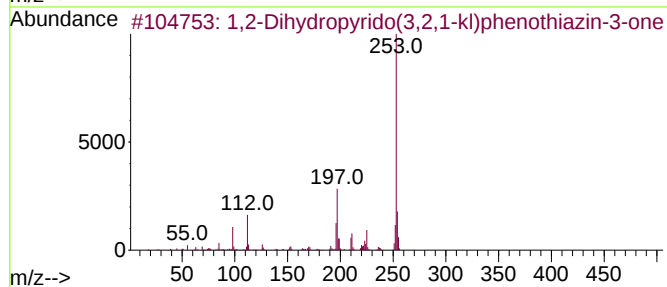
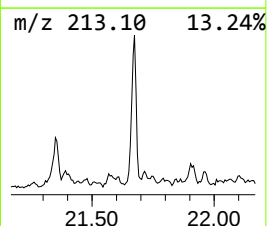
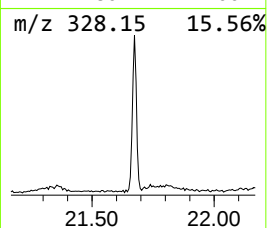
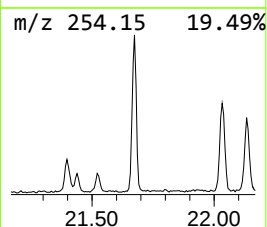
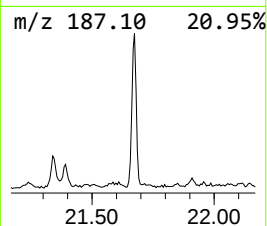
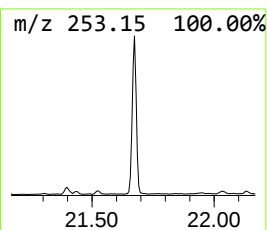
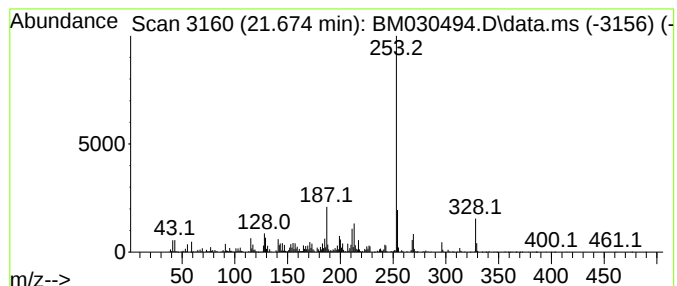
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 1,2-Dihydropyrido(3,2,1-k1)... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.670	4.19 ng/ul	1017030	Chrysene-d12	21.356

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,2-Dihydropyrido(3,2,1-k1)pheno...	253	C15H11NOS	069513-42-4	50
2			Imidazole, 2-iodo-5-methyl-4-nitro-	253	C4H4IN3O2	149510-86-1	43
3			Succinic acid, di(2-phenylphenyl)...	422	C28H22O4	1000330-01-1	43
4			7-(2-Hydroxyphenyl)(7H)triazolo[...	253	C12H7N5O2	166766-15-0	43
5			3,6,6-Trimethyl-1-0-tolyl-1,5,6,...	268	C17H20N2O	1000295-04-0	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061721\
Data File : BM030494.D
Acq On : 17 Jun 2021 14:28
Operator : CG/JU
Sample : M2596-16DL 2X
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM060621.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2H-Pyran, tetra...	3.840	4.4	ng/ul	310948	1	7.828	1399710	20.0
Phenanthrene, 2...	18.060	4.0	ng/ul	511659	4	17.186	2541800	20.0
4H-Cyclopenta[d...	18.260	5.0	ng/ul	637173	4	17.186	2541800	20.0
Anthracene, 2-m...	18.290	2.0	ng/ul	260996	4	17.186	2541800	20.0
1,2-Benzenedica...	18.860	8.9	ng/ul	1132550	4	17.186	2541800	20.0
trans-1,2-Bis(m...	20.540	3.5	ng/ul	844423	5	21.356	4859960	20.0
11H-Benzo[a]flu...	21.130	2.3	ng/ul	545542	5	21.356	4859960	20.0
1,2-Dihydropyri...	21.670	4.2	ng/ul	1017030	5	21.356	4859960	20.0