

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM060424\
 Data File : BM045821.D
 Acq On : 04 Jun 2024 17:53
 Operator : MA/JU
 Sample : P2589-10
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BH6C4

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-BM053124.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BM045821.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.051	9	11	21	rVB	61240	99379	1.46%	0.101%
2	3.475	81	83	92	rVV	242229	546995	8.05%	0.556%
3	3.551	92	96	104	rVV	253167	442114	6.50%	0.450%
4	3.616	105	107	117	rVB3	46063	81291	1.20%	0.083%
5	6.716	628	634	648	rBV	1109524	1990726	29.28%	2.025%
6	6.834	648	654	668	rVB	857883	1453744	21.38%	1.479%
7	7.028	680	687	696	rBV	1195075	1878308	27.63%	1.911%
8	7.104	696	700	710	rVB6	24805	58507	0.86%	0.060%
9	7.463	755	761	770	rVB	1111799	1753003	25.79%	1.783%
10	8.234	886	892	901	rBV	1063460	1758038	25.86%	1.788%
11	8.628	953	959	969	rBV	1190634	1986731	29.22%	2.021%
12	9.345	1071	1081	1096	rBV	930877	1610095	23.68%	1.638%
13	9.892	1167	1174	1189	rVV	1364974	2271428	33.41%	2.311%
14	10.233	1224	1232	1238	rBV	1602831	2553037	37.55%	2.597%
15	10.286	1238	1241	1251	rVB3	28547	58387	0.86%	0.059%
16	10.422	1257	1264	1278	rBV	635691	1329238	19.55%	1.352%
17	11.692	1472	1480	1486	rVV2	42383	70419	1.04%	0.072%
18	11.833	1500	1504	1512	rVV3	35440	60425	0.89%	0.061%
19	12.957	1685	1695	1701	rBV2	44166	72100	1.06%	0.073%
20	13.063	1708	1713	1723	rBV	35972	79266	1.17%	0.081%
21	13.557	1789	1797	1805	rBV	2251787	3271346	48.12%	3.328%
22	13.798	1831	1838	1851	rBV	2604089	3960143	58.25%	4.028%
23	14.033	1873	1878	1882	rBV	38185	55910	0.82%	0.057%
24	14.110	1884	1891	1898	rBV	2194792	3148183	46.31%	3.202%
25	14.174	1898	1902	1906	rVV	46943	75919	1.12%	0.077%
26	14.233	1909	1912	1923	rVB10	29317	56704	0.83%	0.058%
27	14.345	1923	1931	1935	rBV2	94916	147129	2.16%	0.150%
28	14.404	1935	1941	1952	rVV	997286	1496976	22.02%	1.523%
29	14.515	1956	1960	1962	rVV2	53729	88424	1.30%	0.090%
30	14.539	1962	1964	1967	rVV3	52239	72551	1.07%	0.074%
31	14.992	2033	2041	2045	rVB2	47359	66379	0.98%	0.068%
32	15.110	2054	2061	2067	rBV	3337692	4616340	67.90%	4.696%
33	15.245	2079	2084	2089	rBV	1013918	1294485	19.04%	1.317%
34	15.615	2142	2147	2154	rVB4	30224	67549	0.99%	0.069%
35	15.851	2182	2187	2190	rBV3	74621	100740	1.48%	0.102%
36	16.168	2235	2241	2246	rBV4	30409	64256	0.95%	0.065%

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Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM053124.MA.M
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37	16.521	2296	2301	2307	rVB2	78482	125875	1.85%	0.128%
38	16.639	2319	2321	2324	rVV2	51047	64023	0.94%	0.065%
39	16.674	2324	2327	2337	rVB2	73763	137065	2.02%	0.139%
40	16.851	2352	2357	2361	rBV	2643137	3645149	53.62%	3.708%
41	16.892	2361	2364	2369	rVB	1168854	1566177	23.04%	1.593%
42	16.951	2369	2374	2379	rBV	3327626	4687630	68.95%	4.768%
43	17.068	2388	2394	2398	rBV5	29754	75227	1.11%	0.077%
44	17.174	2408	2412	2417	rVB3	62607	83454	1.23%	0.085%
45	17.274	2422	2429	2436	rBV2	135819	264971	3.90%	0.270%
46	17.374	2443	2446	2452	rBV3	50742	90881	1.34%	0.092%
47	17.439	2453	2457	2467	rVB5	52463	98068	1.44%	0.100%
48	17.656	2490	2494	2501	rBV5	32652	89870	1.32%	0.091%
49	17.727	2501	2506	2510	rVV	307470	400143	5.89%	0.407%
50	17.780	2510	2515	2523	rVV2	544004	1026877	15.11%	1.045%
51	17.851	2523	2527	2533	rVV2	221655	354035	5.21%	0.360%
52	17.915	2533	2538	2542	rVV2	367794	614672	9.04%	0.625%
53	17.956	2542	2545	2554	rVB	213513	289130	4.25%	0.294%
54	18.056	2557	2562	2565	rBV3	74873	111453	1.64%	0.113%
55	18.233	2588	2592	2596	rBV	169313	233961	3.44%	0.238%
56	18.274	2596	2599	2605	rVB	240058	313292	4.61%	0.319%
57	18.480	2631	2634	2639	rVV	97732	136329	2.01%	0.139%
58	18.533	2639	2643	2646	rVV	109807	133480	1.96%	0.136%
59	18.574	2646	2650	2654	rVB	75928	90655	1.33%	0.092%
60	18.656	2659	2664	2669	rVV2	240058	444409	6.54%	0.452%
61	18.703	2669	2672	2677	rVV4	170891	299220	4.40%	0.304%
62	18.745	2677	2679	2683	rVB	82282	92066	1.35%	0.094%
63	18.798	2683	2688	2696	rBV2	211439	431359	6.35%	0.439%
64	18.915	2700	2708	2716	rVB	3152918	3994877	58.76%	4.064%
65	18.992	2718	2721	2724	rVV	57533	75728	1.11%	0.077%
66	19.021	2724	2726	2731	rVB2	64852	80021	1.18%	0.081%
67	19.074	2731	2735	2738	rBV2	108281	141667	2.08%	0.144%
68	19.162	2747	2750	2754	rVB2	98369	131310	1.93%	0.134%
69	19.250	2759	2765	2768	rVV	4940260	6798254	100.00%	6.915%
70	19.280	2768	2770	2779	rVV	3101147	3383592	49.77%	3.442%
71	19.356	2780	2783	2788	rVB3	147825	221320	3.26%	0.225%
72	19.409	2788	2792	2796	rBV	156342	172847	2.54%	0.176%
73	19.462	2798	2801	2808	rVB	140393	191742	2.82%	0.195%
74	19.609	2821	2826	2833	rBV4	236891	525493	7.73%	0.535%
75	19.745	2845	2849	2851	rBV3	208589	288078	4.24%	0.293%
76	19.780	2852	2855	2859	rVB2	295718	402966	5.93%	0.410%

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 Title : SVOA CALIBRATION

77	19.821	2859	2862	2866	rVB	153640	152936	2.25%	0.156%
78	19.880	2869	2872	2876	rBV	167011	206548	3.04%	0.210%
79	19.939	2877	2882	2886	rBV2	344805	507781	7.47%	0.517%
80	20.074	2902	2905	2909	rBV	199445	241797	3.56%	0.246%
81	20.121	2910	2913	2918	rVB2	181253	250638	3.69%	0.255%
82	20.209	2925	2928	2932	rBV3	165988	193060	2.84%	0.196%
83	20.315	2943	2946	2949	rBV	184240	199560	2.94%	0.203%
84	20.356	2949	2953	2959	rVV	312695	408242	6.01%	0.415%
85	20.533	2979	2983	2989	rVB	325778	441487	6.49%	0.449%
86	20.692	3006	3010	3014	rVB2	230815	324795	4.78%	0.330%
87	20.733	3014	3017	3020	rVV	235390	280474	4.13%	0.285%
88	20.780	3020	3025	3030	rVV3	680041	1165493	17.14%	1.186%
89	20.827	3031	3033	3045	rVB2	319486	565506	8.32%	0.575%
90	21.003	3059	3063	3066	rBV	1041172	1201497	17.67%	1.222%
91	21.062	3066	3073	3077	rVV3	2997654	5825383	85.69%	5.926%
92	21.103	3077	3080	3089	rVB	1376229	1912654	28.13%	1.946%
93	21.768	3188	3193	3198	rBV2	337714	544279	8.01%	0.554%
94	22.350	3288	3292	3300	rVB	354044	595686	8.76%	0.606%
95	22.933	3385	3391	3398	rBV	837310	2286311	33.63%	2.326%
96	23.538	3488	3494	3498	rBV	467832	927847	13.65%	0.944%
97	23.603	3498	3505	3510	rVV2	1746183	3715742	54.66%	3.780%
98	23.656	3510	3514	3526	rVB	649503	1359032	19.99%	1.382%
99	23.785	3528	3536	3543	rBV	1533178	3469267	51.03%	3.529%
100	24.685	3685	3689	3699	rVB	224465	485643	7.14%	0.494%

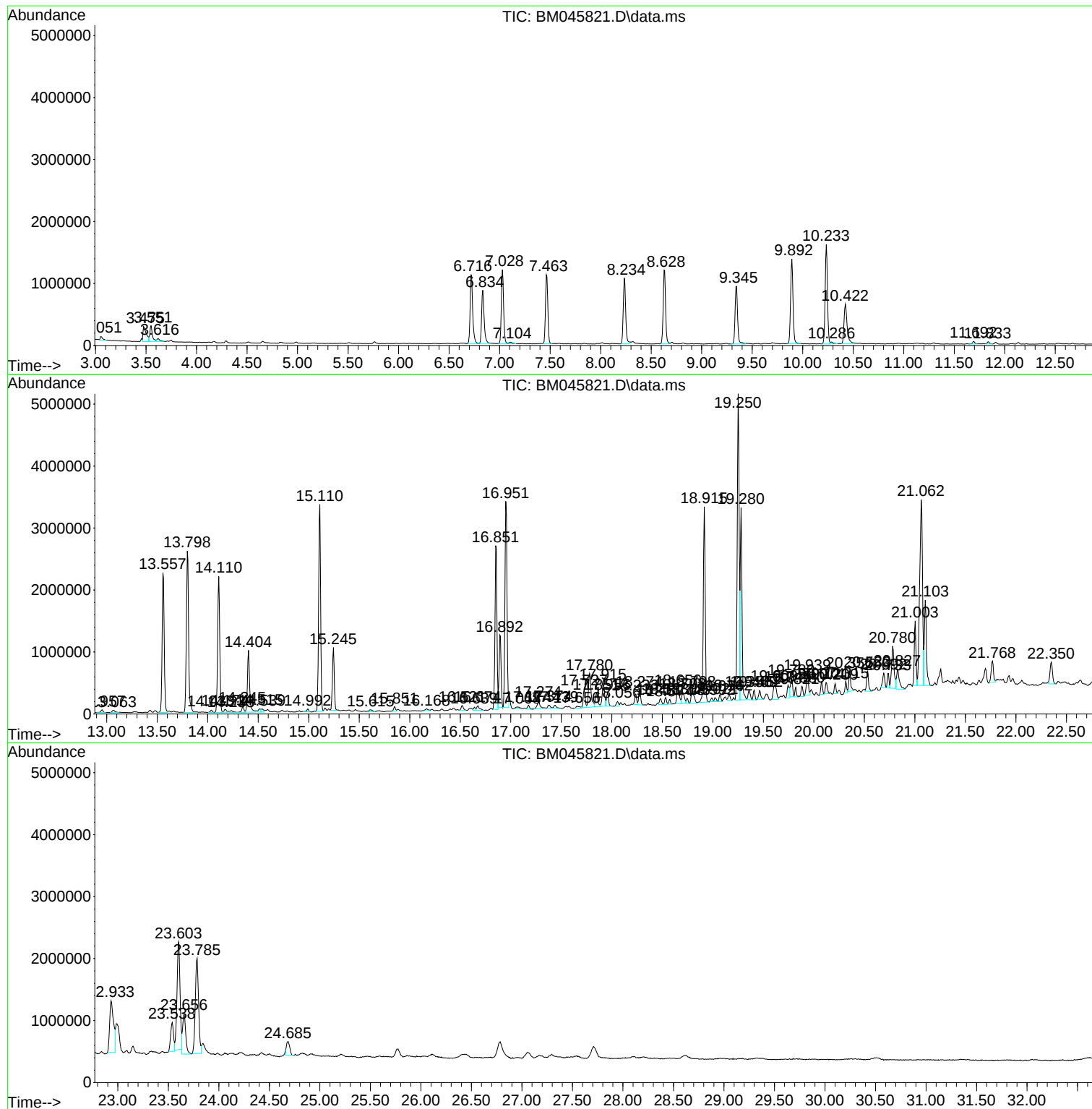
Sum of corrected areas: 98305289

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM053124.MA.M
Quant Title : SVOA CALIBRATION

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



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Instrument :
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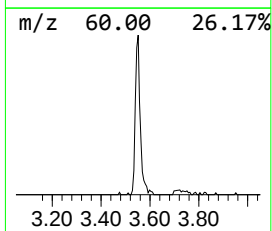
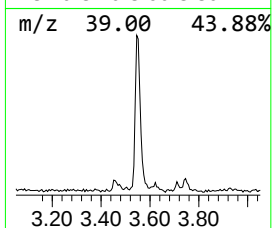
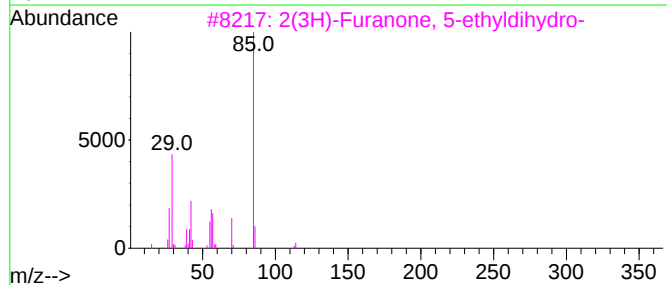
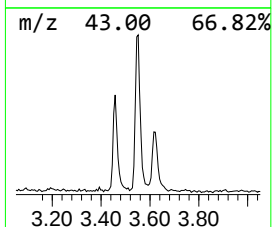
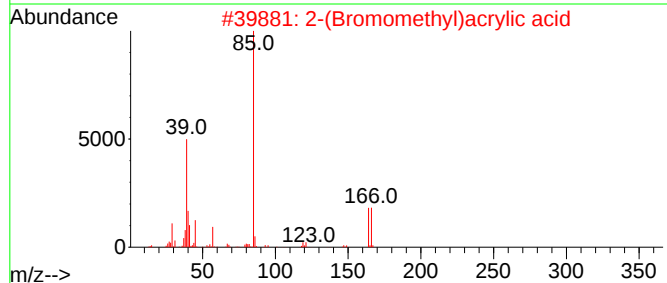
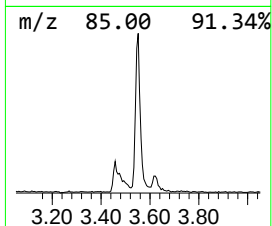
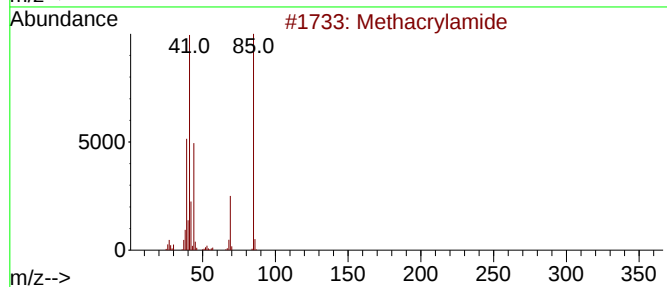
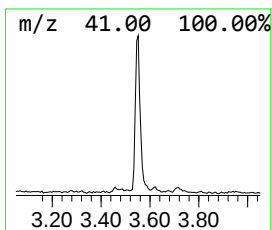
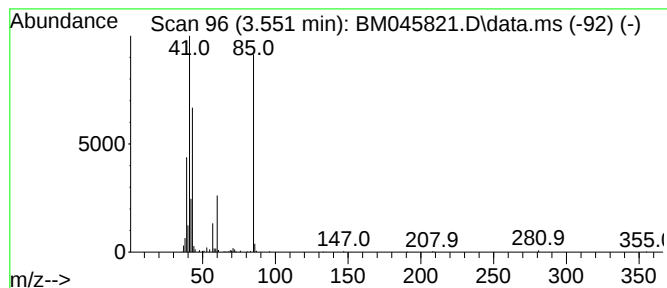
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM053124.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Methacrylamide Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.551	5.04 ng/ul	442114	1,4-Dichlorobenzene-d4	7.469

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methacrylamide	85	C4H7NO	000079-39-0	53
2			2-(Bromomethyl)acrylic acid	164	C4H5BrO2	072707-66-5	40
3			2(3H)-Furanone, 5-ethyldihydro-	114	C6H10O2	000695-06-7	25
4			Sulfurous acid, dodecyl hexyl ester	334	C18H38O3S	1000309-13-4	17
5			Heptane, 4,4-dimethyl-	128	C9H20	001068-19-5	12



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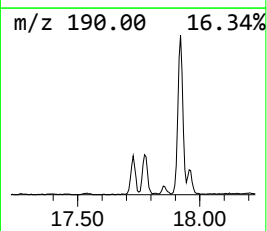
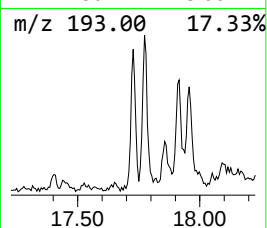
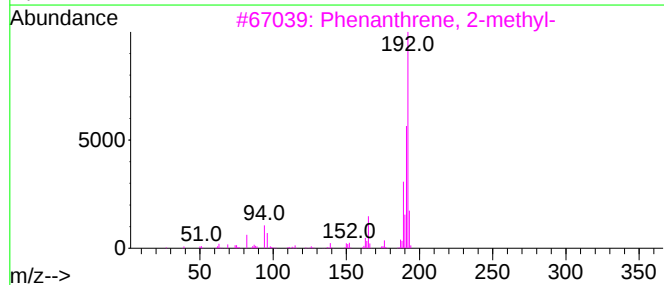
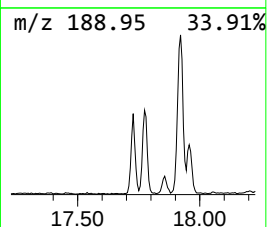
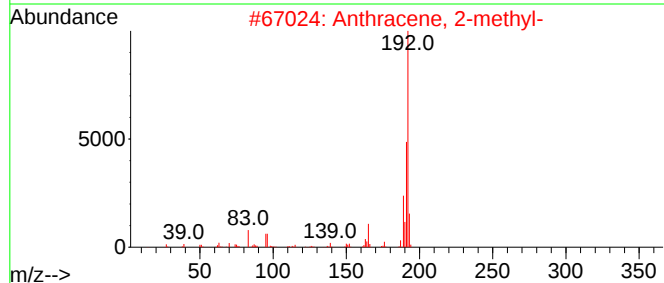
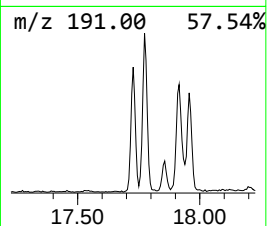
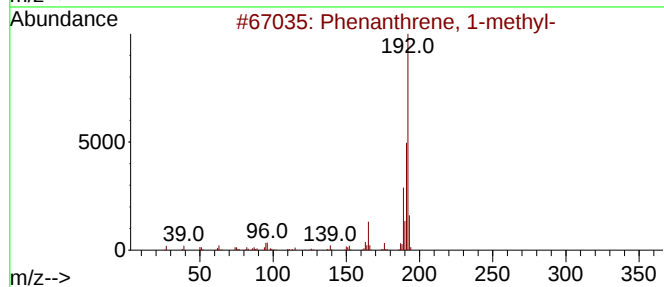
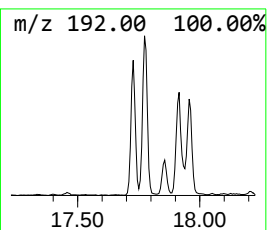
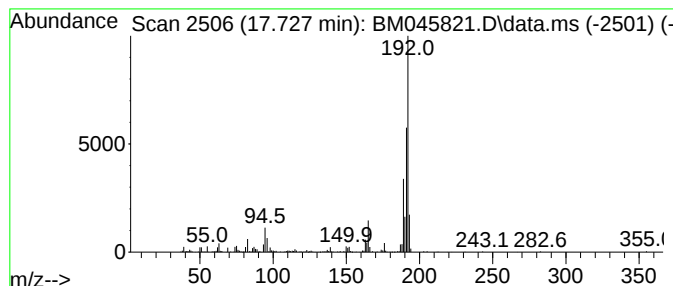
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM053124.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Phenanthrene, 1-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.727	2.20 ng/ul	400143	Phenanthrene-d10	16.851

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	97
2		Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
3		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
4		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95
5		1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	95



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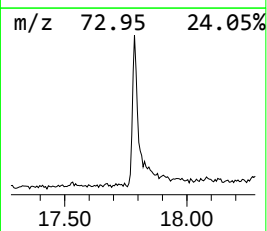
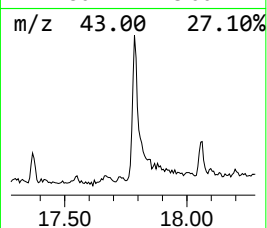
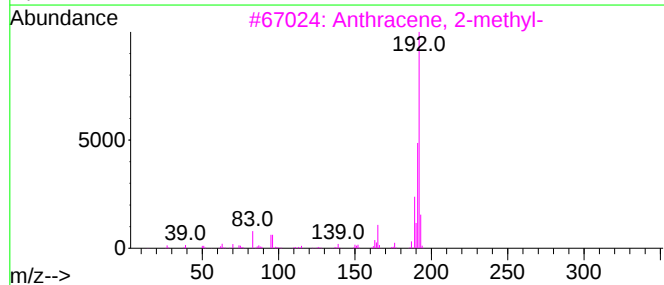
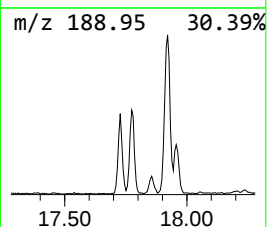
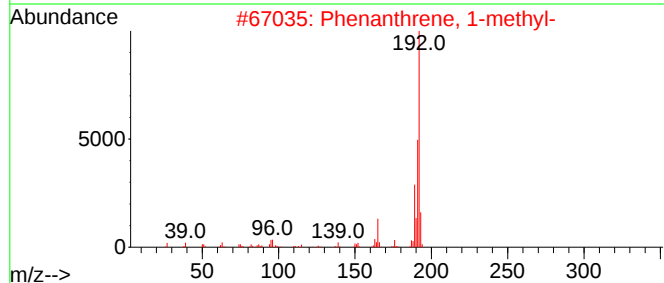
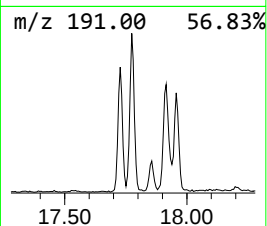
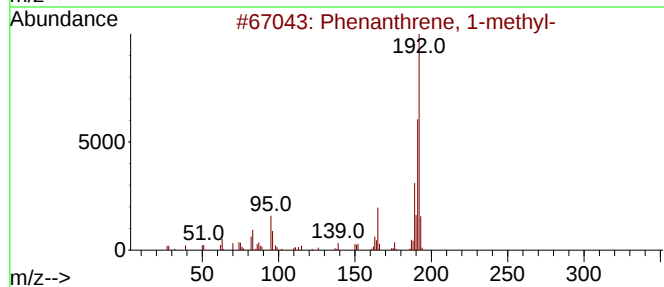
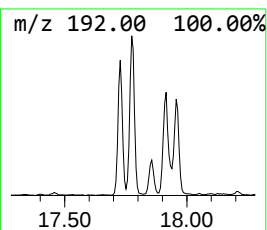
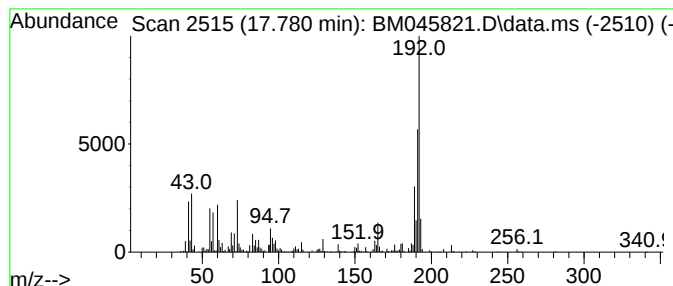
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM053124.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Anthracene, 2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.780	5.63 ng/ul	1026880	Phenanthrene-d10	16.851

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	98
2			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	97
3			Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
4			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
5			1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	95



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Data File : BM045821.D
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Operator : MA/JU
Sample : P2589-10
Misc :
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Instrument :
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ClientSampleId :
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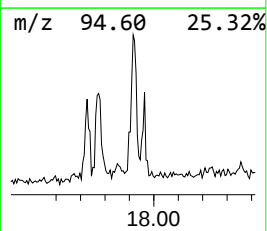
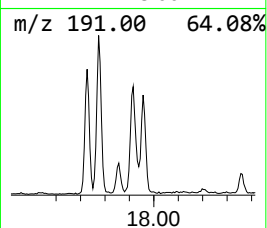
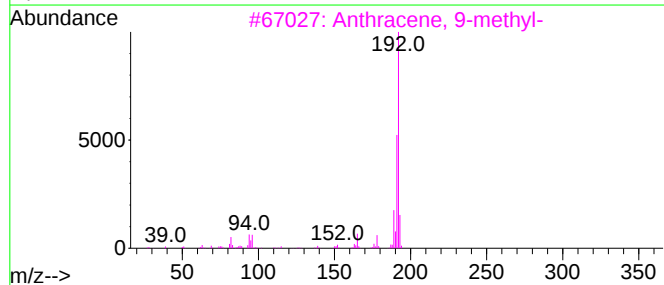
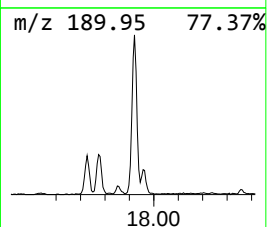
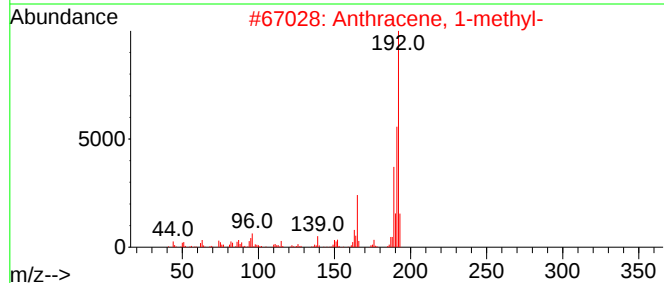
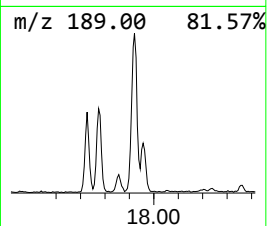
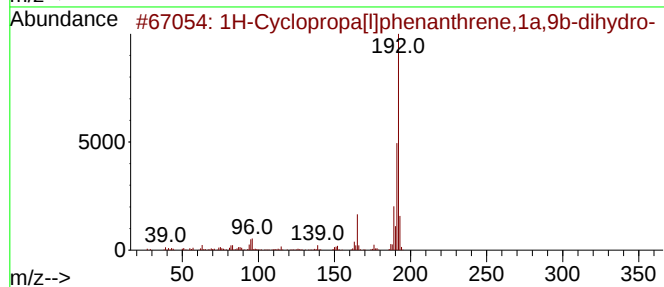
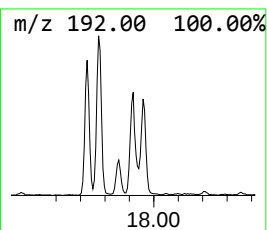
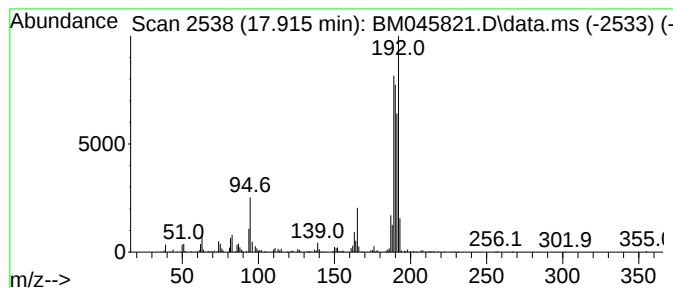
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 unknown-01 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.915	3.37 ng/ul	614672	Phenanthrene-d10	16.851

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	46
2			Anthracene, 1-methyl-	192	C15H12	000610-48-0	46
3			Anthracene, 9-methyl-	192	C15H12	000779-02-2	45
4			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	43
5			Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM060424\
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ClientSampleId :
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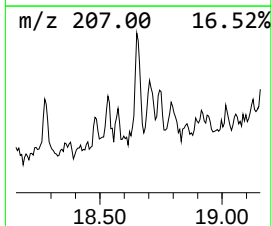
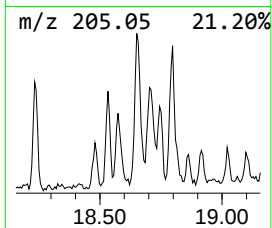
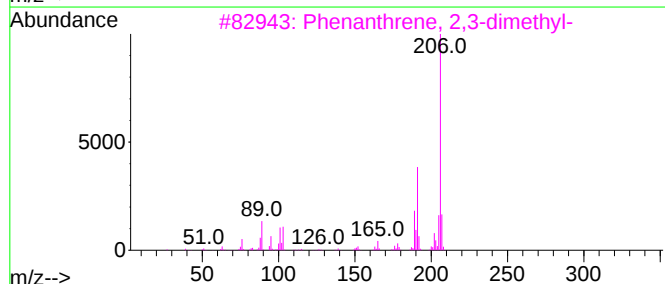
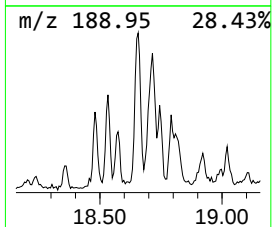
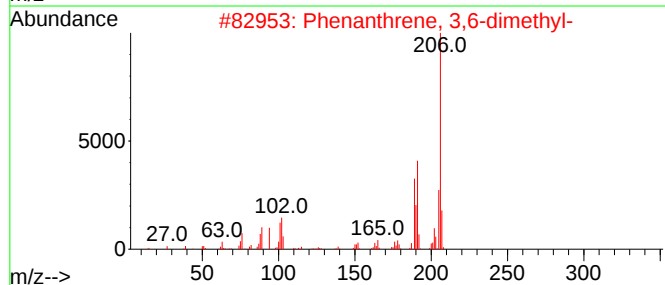
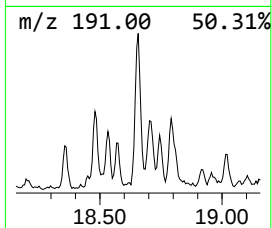
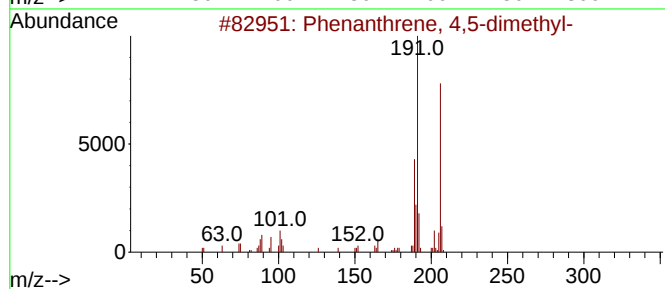
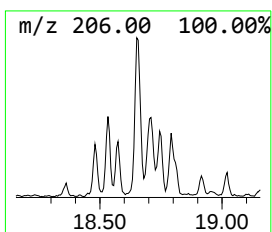
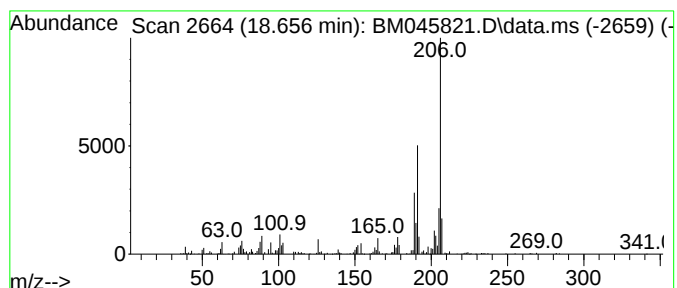
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Phenanthrene, 4,5-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.656	2.44 ng/ul	444409	Phenanthrene-d10	16.851

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 4,5-dimethyl-	206	C16H14	003674-69-9	94
2			Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	93
3			Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	93
4			9,10-Dimethylanthracene	206	C16H14	000781-43-1	92
5			9,10-Dimethylanthracene	206	C16H14	000781-43-1	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM060424\
Data File : BM045821.D
Acq On : 04 Jun 2024 17:53
Operator : MA/JU
Sample : P2589-10
Misc :
ALS Vial : 11 Sample Multiplier: 1

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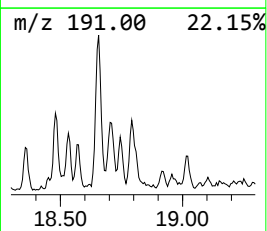
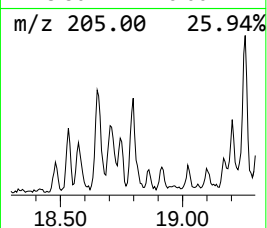
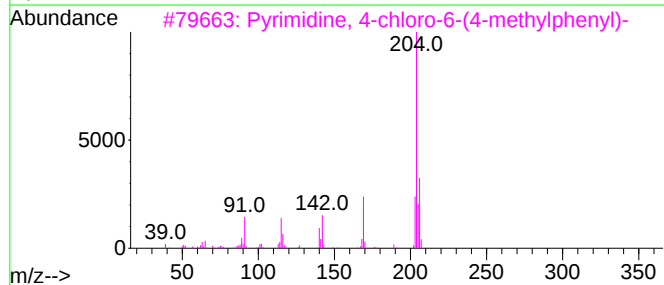
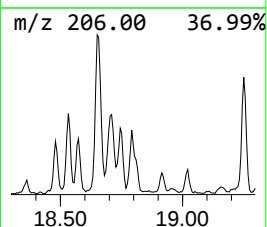
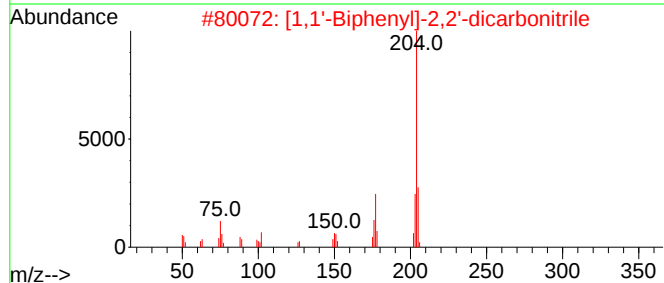
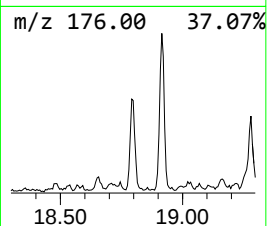
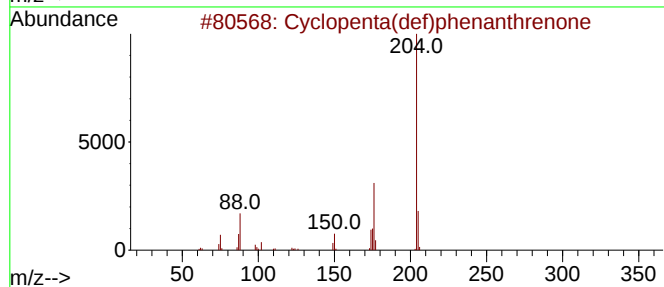
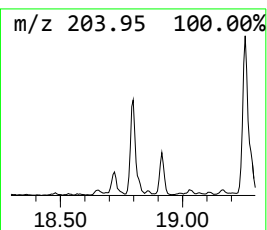
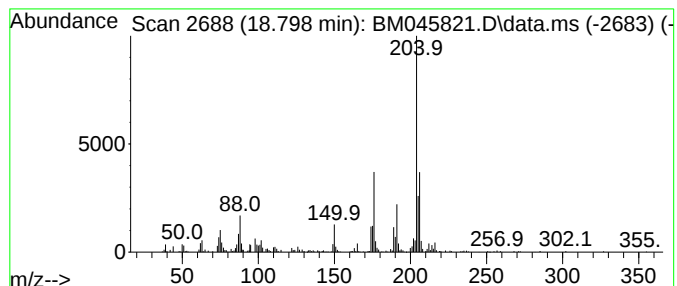
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Cyclopenta(def)phenanthrene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.797	2.37 ng/ul	431359	Phenanthrene-d10	16.851

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopenta(def)phenanthrene	204	C15H8O	005737-13-3	93
2			[1,1'-Biphenyl]-2,2'-dicarbonitrile	204	C14H8N2	004341-02-0	49
3			Pyrimidine, 4-chloro-6-(4-methyl...	204	C11H9ClN2	1000380-55-8	47
4			Benzene, 1-chloro-4-phenoxy-	204	C12H9ClO	007005-72-3	46
5			1,2,4,8-Tetramethylbicyclo[6.3.0...	204	C15H24	137235-51-9	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM060424\
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Sample : P2589-10
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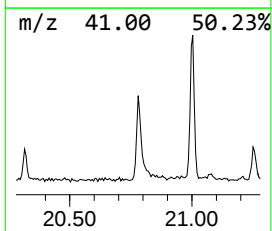
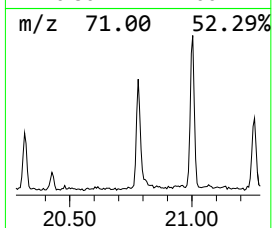
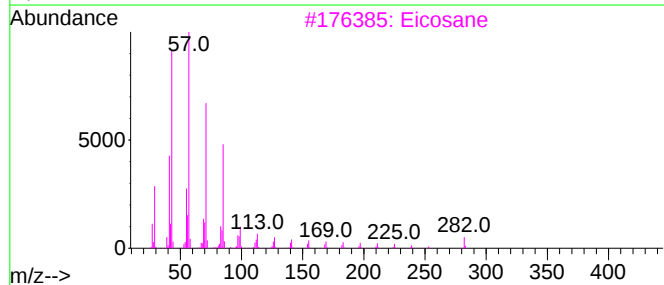
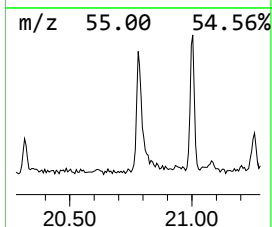
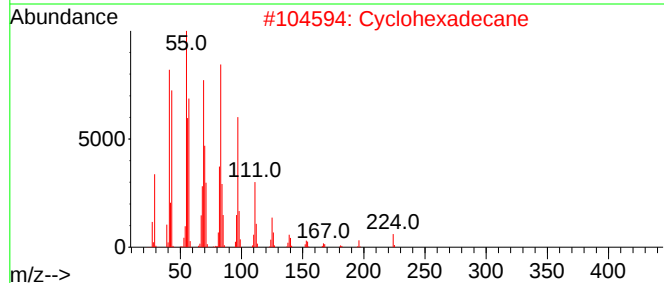
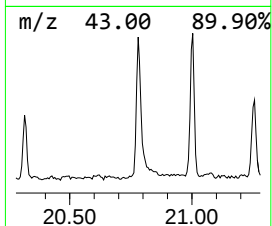
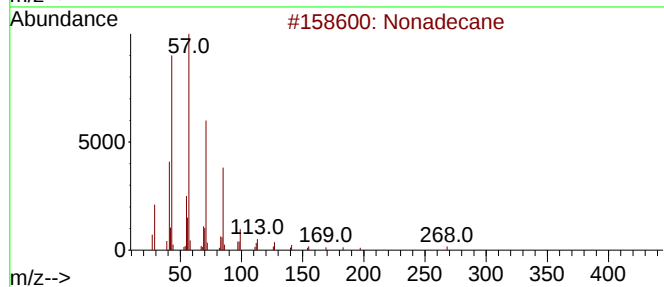
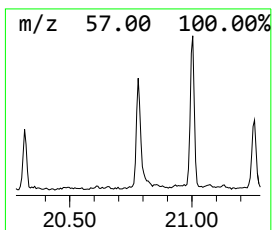
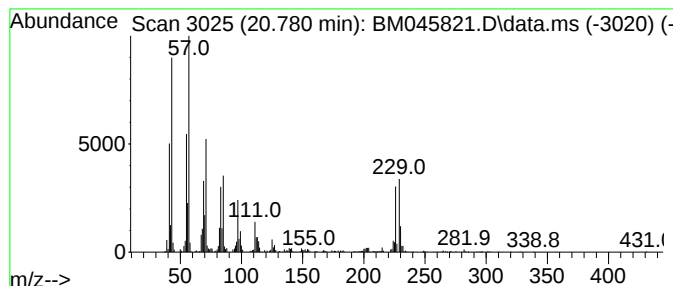
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
20.780	4.00 ng/ul	1165490	Chrysene-d12	21.062	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonadecane	268	C19H40	000629-92-5	83
2	Cyclohexadecane	224	C16H32	000295-65-8	64
3	Eicosane	282	C20H42	000112-95-8	60
4	Hexadecane	226	C16H34	000544-76-3	60
5	Carbonic acid, eicosyl vinyl ester	368	C23H44O3	1000382-54-3	60



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM060424\
Data File : BM045821.D
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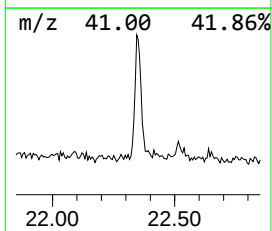
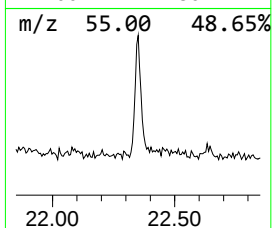
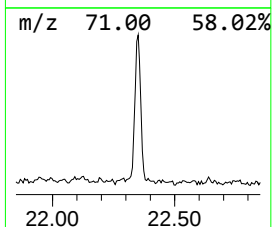
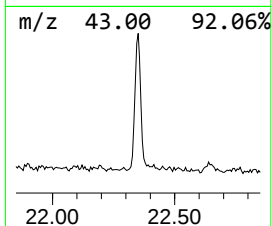
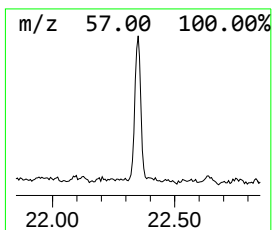
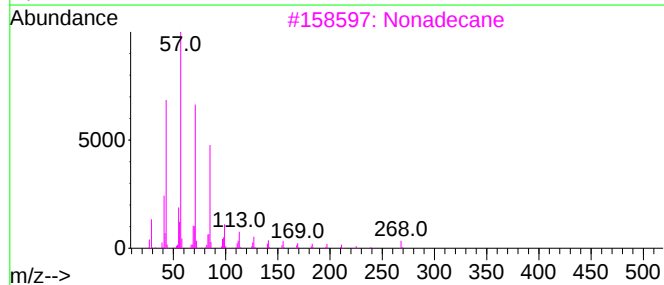
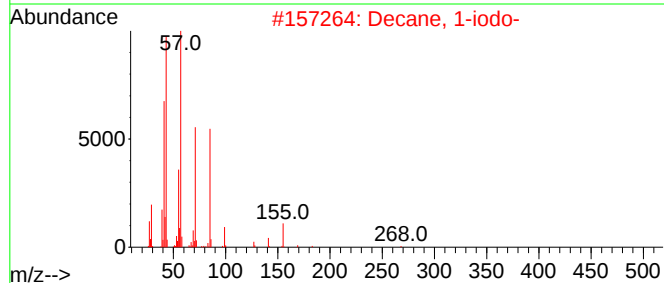
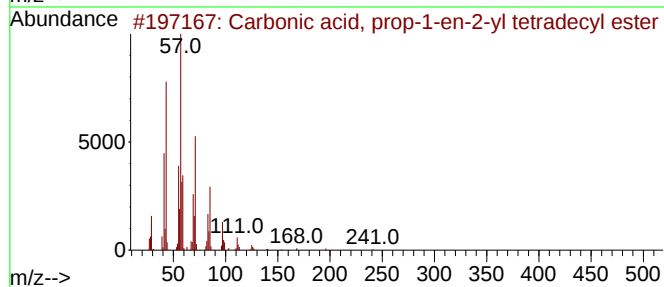
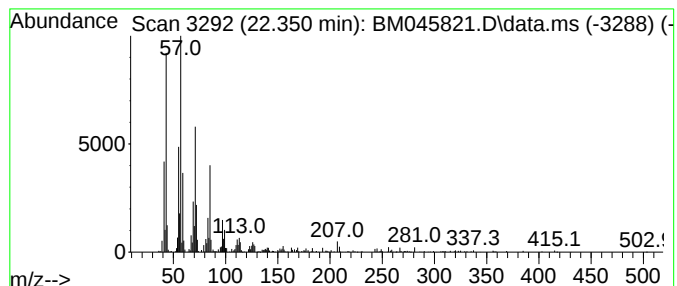
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Carbonic acid, prop-1-en-2-... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.350	2.05 ng/ul	595686	Chrysene-d12	21.062

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Carbonic acid, prop-1-en-2-yl te...	298	C18H34O3	1000382-90-9	68
2			Decane, 1-iodo-	268	C10H21I	002050-77-3	64
3			Nonadecane	268	C19H40	000629-92-5	64
4			Octacosane	394	C28H58	000630-02-4	62
5			Nonadecane	268	C19H40	000629-92-5	62



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM060424\
Data File : BM045821.D
Acq On : 04 Jun 2024 17:53
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ALS Vial : 11 Sample Multiplier: 1

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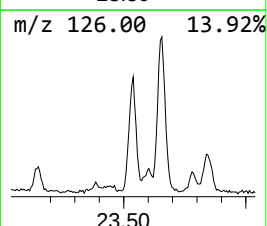
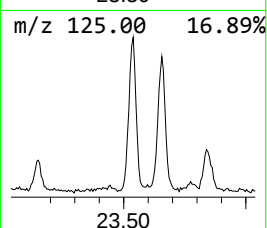
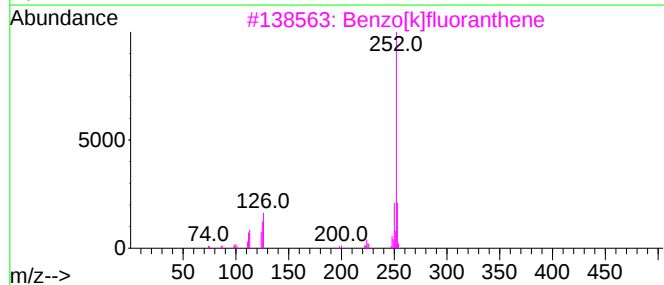
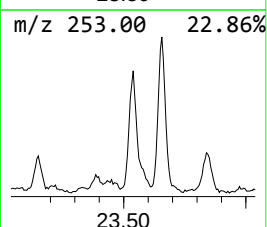
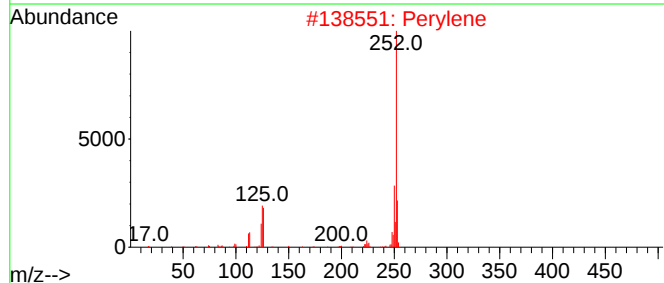
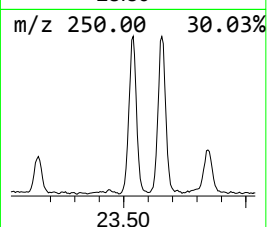
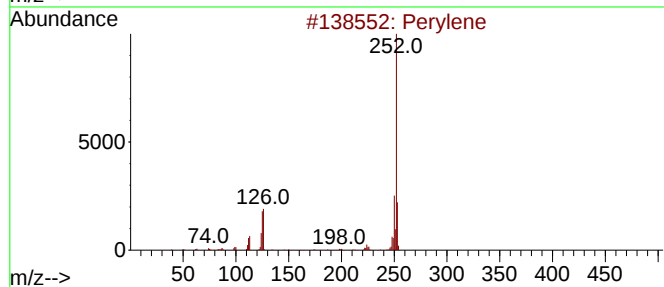
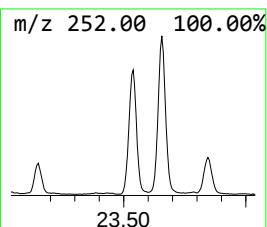
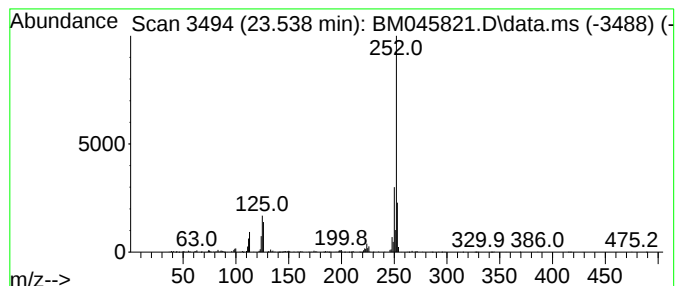
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Perylene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.538	5.35 ng/ul	927847	Perylene-d12	23.785

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Perylene	252	C20H12	000198-55-0	99
2			Perylene	252	C20H12	000198-55-0	98
3			Benzo[k]fluoranthene	252	C20H12	000207-08-9	98
4			Benzo[e]pyrene	252	C20H12	000192-97-2	98
5			Benzo[b]fluoranthene	252	C20H12	000205-99-2	98



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM060424\
Data File : BM045821.D
Acq On : 04 Jun 2024 17:53
Operator : MA/JU
Sample : P2589-10
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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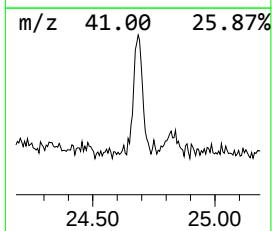
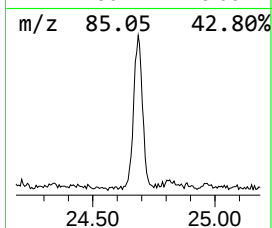
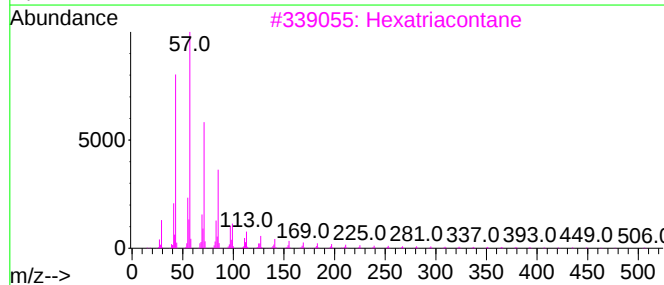
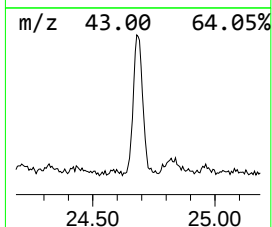
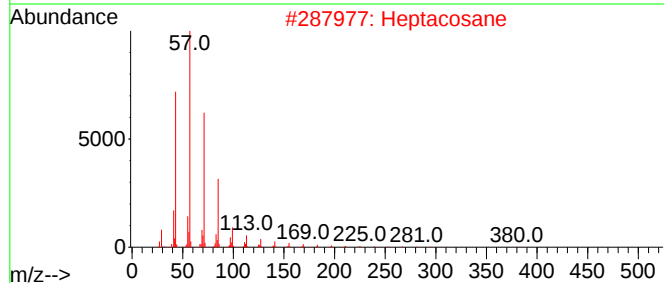
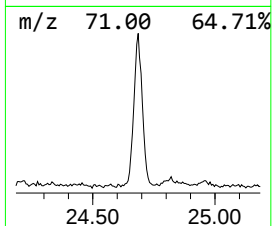
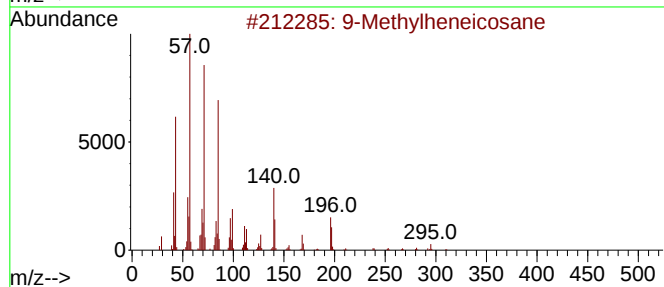
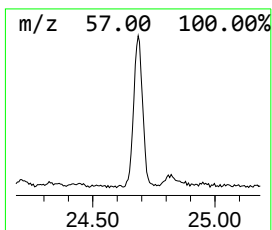
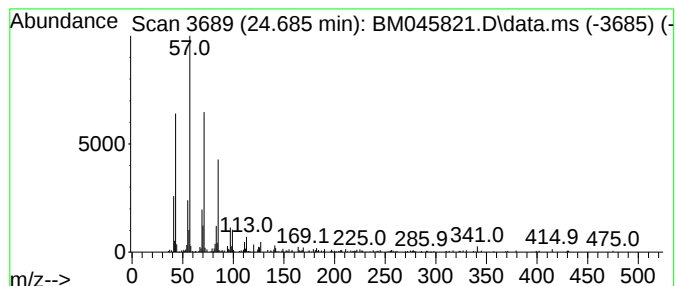
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM053124.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.685	2.80 ng/ul	485643	Perylene-d12	23.785

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		9-Methylheneicosane	310	C22H46	070475-51-3	86
2		Heptacosane	380	C27H56	000593-49-7	80
3		Hexatriacontane	507	C36H74	000630-06-8	64
4		Sulfurous acid, 2-ethylhexyl oct...	446	C26H54O3S	1000309-20-1	52
5		Tetratetracontane	619	C44H90	007098-22-8	52



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM060424\
Data File : BM045821.D
Acq On : 04 Jun 2024 17:53
Operator : MA/JU
Sample : P2589-10
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BH6C4

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM053124.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
Methacrylamide	3.551	5.0	ng/ul	442114	1	7.469	1753000	20.0
Phenanthrene, 1...	17.727	2.2	ng/ul	400143	4	16.851	3645150	20.0
Anthracene, 2-m...	17.780	5.6	ng/ul	1026880	4	16.851	3645150	20.0
unknown-01	17.915	3.4	ng/ul	614672	4	16.851	3645150	20.0
Phenanthrene, 4...	18.656	2.4	ng/ul	444409	4	16.851	3645150	20.0
Cyclopenta(def)...	18.797	2.4	ng/ul	431359	4	16.851	3645150	20.0
(DEL) Alkane: S...	20.780	4.0	ng/ul	1165490	5	21.062	5825380	20.0
Carbonic acid, ...	22.350	2.0	ng/ul	595686	5	21.062	5825380	20.0
Perylene	23.538	5.3	ng/ul	927847	6	23.785	3469270	20.0
(DEL) Alkane: S...	24.685	2.8	ng/ul	485643	6	23.785	3469270	20.0