

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100724\
 Data File : BM047933.D
 Acq On : 09 Oct 2024 04:56
 Operator : RC/JU
 Sample : P4183-05
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4EM1

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-BM092724.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BM047933.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	41	44	53	rVB5	26017	43181	0.59%	0.051%
2	3.681	114	118	123	rBV2	19803	29864	0.41%	0.035%
3	3.763	128	132	138	rVB	45683	58755	0.80%	0.070%
4	3.987	167	170	175	rVB	25210	30885	0.42%	0.037%
5	4.904	321	326	333	rVB	45215	70104	0.96%	0.083%
6	6.051	517	521	526	rVB4	22780	32430	0.44%	0.039%
7	6.787	640	646	650	rBV2	23278	39953	0.55%	0.047%
8	6.963	667	676	691	rVB	513135	951537	12.99%	1.131%
9	7.122	697	703	714	rVB	445238	690669	9.43%	0.821%
10	7.328	730	738	747	rBV	557804	887312	12.11%	1.054%
11	7.792	810	817	825	rBV	1049221	1641122	22.40%	1.950%
12	7.863	825	829	835	rVB5	16844	30229	0.41%	0.036%
13	8.498	930	937	946	rBV	558581	907157	12.38%	1.078%
14	8.610	951	956	963	rVB	82111	140147	1.91%	0.167%
15	8.945	1006	1013	1022	rBV	488341	804761	10.99%	0.956%
16	9.510	1100	1109	1118	rBV2	49974	91107	1.24%	0.108%
17	9.669	1129	1136	1143	rBV	375922	640428	8.74%	0.761%
18	9.822	1156	1162	1166	rBV2	20314	41316	0.56%	0.049%
19	10.028	1188	1197	1200	rBV8	12991	29290	0.40%	0.035%
20	10.210	1219	1228	1237	rBV	602775	1041364	14.22%	1.237%
21	10.586	1284	1292	1298	rBV	1345091	2297447	31.36%	2.730%
22	10.633	1298	1300	1307	rVV	42402	63542	0.87%	0.075%
23	10.722	1307	1315	1334	rVB	121467	287177	3.92%	0.341%
24	11.122	1378	1383	1390	rBV7	23259	41770	0.57%	0.050%
25	11.875	1507	1511	1516	rBV6	15483	26974	0.37%	0.032%
26	12.251	1568	1575	1580	rBV3	22522	39242	0.54%	0.047%
27	12.363	1585	1594	1599	rBV5	12365	31410	0.43%	0.037%
28	13.257	1743	1746	1752	rVV4	25920	39874	0.54%	0.047%
29	13.604	1795	1805	1810	rBV8	12120	30411	0.42%	0.036%
30	13.745	1826	1829	1835	rBV5	27033	42691	0.58%	0.051%
31	13.833	1835	1844	1850	rBV	970263	1367643	18.67%	1.625%
32	14.121	1882	1893	1903	rBV	1143166	1953390	26.67%	2.321%
33	14.427	1936	1945	1951	rBV	1991889	2866089	39.13%	3.405%
34	14.492	1951	1956	1963	rVB	65081	103651	1.41%	0.123%
35	14.551	1963	1966	1971	rVB6	25154	35509	0.48%	0.042%
36	14.627	1971	1979	1993	rBV	371628	696595	9.51%	0.828%

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Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
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37	14.827	2008	2013	2020	rVB	48229	90408	1.23%	0.107%
38	14.904	2020	2026	2032	rVB5	19723	32008	0.44%	0.038%
39	15.227	2074	2081	2087	rBV7	19154	48283	0.66%	0.057%
40	15.286	2087	2091	2097	rVB2	17463	34728	0.47%	0.041%
41	15.421	2107	2114	2119	rBV	1483506	2183183	29.80%	2.594%
42	15.474	2119	2123	2127	rVV2	66902	100401	1.37%	0.119%
43	15.533	2128	2133	2139	rVV	336816	475986	6.50%	0.566%
44	15.680	2154	2158	2163	rVB3	41869	62754	0.86%	0.075%
45	15.780	2169	2175	2185	rVB	615577	873640	11.93%	1.038%
46	15.915	2194	2198	2205	rVB	26789	52764	0.72%	0.063%
47	15.992	2205	2211	2218	rVB10	24833	63443	0.87%	0.075%
48	16.145	2232	2237	2243	rBV4	53662	113455	1.55%	0.135%
49	16.239	2250	2253	2257	rBV2	22673	37611	0.51%	0.045%
50	16.345	2267	2271	2278	rBV	39845	69214	0.94%	0.082%
51	16.474	2291	2293	2298	rVB5	21255	28782	0.39%	0.034%
52	16.574	2306	2310	2315	rBV6	31858	47533	0.65%	0.056%
53	16.704	2330	2332	2336	rVB5	22969	26938	0.37%	0.032%
54	16.821	2348	2352	2359	rVB	81636	124701	1.70%	0.148%
55	16.898	2361	2365	2366	rBV4	22752	28824	0.39%	0.034%
56	16.927	2366	2370	2376	rVV3	81883	127643	1.74%	0.152%
57	16.986	2376	2380	2385	rVB	76090	102939	1.41%	0.122%
58	17.068	2390	2394	2401	rVB3	82732	134005	1.83%	0.159%
59	17.168	2401	2411	2414	rBV	2230624	3351919	45.76%	3.983%
60	17.210	2414	2418	2423	rVV	1497723	2117870	28.91%	2.516%
61	17.262	2423	2427	2431	rVV	1628209	2364949	32.28%	2.810%
62	17.298	2431	2433	2438	rVV	275575	335712	4.58%	0.399%
63	17.357	2438	2443	2449	rVV2	72539	123649	1.69%	0.147%
64	17.568	2470	2479	2484	rBV2	177798	368871	5.04%	0.438%
65	17.674	2493	2497	2498	rBV3	36185	51643	0.70%	0.061%
66	17.745	2506	2509	2516	rVB4	63375	105601	1.44%	0.125%
67	17.945	2540	2543	2549	rVB7	45750	72932	1.00%	0.087%
68	18.004	2550	2553	2555	rBV2	37258	48526	0.66%	0.058%
69	18.062	2555	2563	2566	rBV2	687418	1217424	16.62%	1.447%
70	18.168	2578	2581	2587	rVB	57060	71934	0.98%	0.085%
71	18.233	2587	2592	2596	rBV2	353626	627148	8.56%	0.745%
72	18.268	2596	2598	2606	rVB	139732	211141	2.88%	0.251%
73	18.356	2607	2613	2622	rBV6	103757	275048	3.75%	0.327%
74	18.539	2640	2644	2647	rBV	180916	242722	3.31%	0.288%
75	18.580	2647	2651	2656	rVB2	340421	459165	6.27%	0.546%
76	18.786	2683	2686	2690	rVB2	94064	110548	1.51%	0.131%

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 Title : SVOA CALIBRATION

77	18.962	2712	2716	2718	rBV2	152019	209803	2.86%	0.249%
78	18.992	2718	2721	2724	rVV	117649	131905	1.80%	0.157%
79	19.092	2735	2738	2747	rVB	195978	336737	4.60%	0.400%
80	19.221	2755	2760	2767	rBV	4722940	6209546	84.77%	7.378%
81	19.339	2777	2780	2782	rBV	144415	145179	1.98%	0.172%
82	19.551	2811	2816	2819	rBV2	2408247	3746873	51.15%	4.452%
83	19.580	2819	2821	2829	rVV	2745599	3482411	47.54%	4.138%
84	19.651	2830	2833	2839	rVB2	168227	250516	3.42%	0.298%
85	19.762	2849	2852	2856	rVB	170963	209222	2.86%	0.249%
86	19.903	2872	2876	2883	rBV4	224334	517093	7.06%	0.614%
87	20.074	2901	2905	2908	rBV	505133	627126	8.56%	0.745%
88	20.174	2919	2922	2926	rVB	350983	388609	5.30%	0.462%
89	20.821	3029	3032	3037	rVB	404926	533679	7.29%	0.634%
90	21.039	3066	3069	3071	rBV	295497	374672	5.11%	0.445%
91	21.068	3071	3074	3081	rVB3	1588739	2290545	31.27%	2.722%
92	21.297	3109	3113	3117	rVB	855588	1049325	14.32%	1.247%
93	21.392	3121	3129	3133	rVV4	2924545	7325412	100.00%	8.704%
94	21.433	3133	3136	3149	rVB	2252897	3616353	49.37%	4.297%
95	22.156	3254	3259	3261	rBV2	1262705	2205039	30.10%	2.620%
96	23.121	3419	3423	3435	rVB2	772886	1567386	21.40%	1.862%
97	23.444	3472	3478	3485	rBV	1833935	5000533	68.26%	5.942%
98	23.603	3500	3505	3514	rVB2	1347231	2820926	38.51%	3.352%
99	24.397	3633	3640	3649	rVB	1473025	3634751	49.62%	4.319%
100	25.450	3814	3819	3829	rVB2	896453	2347496	32.05%	2.789%

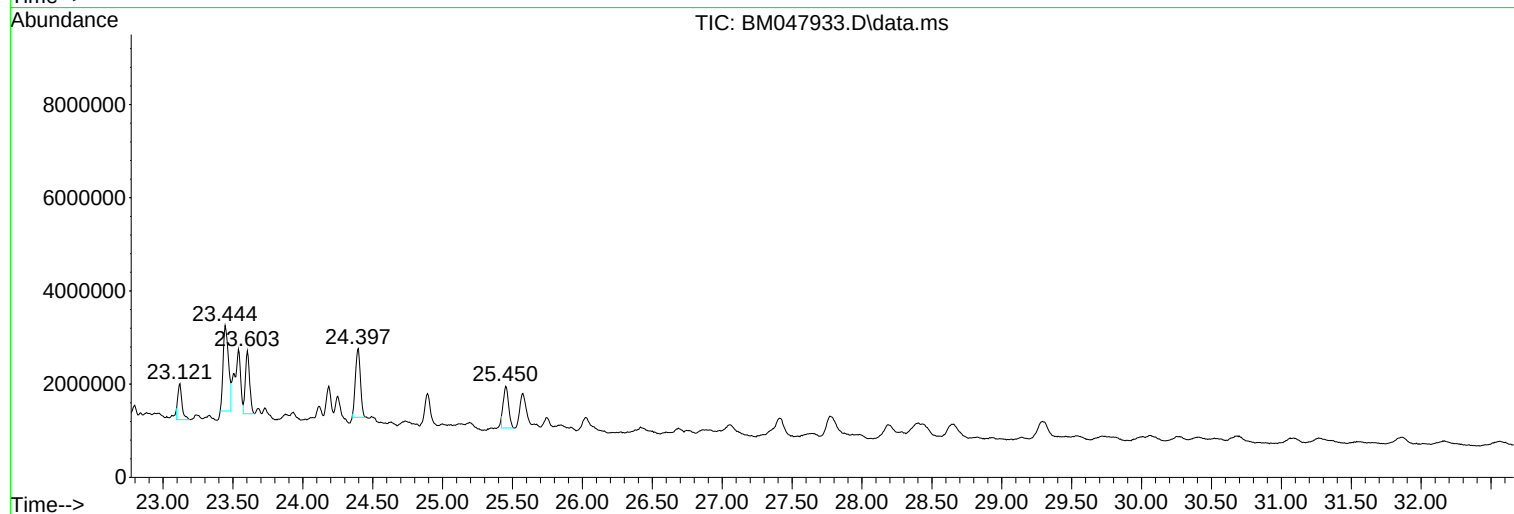
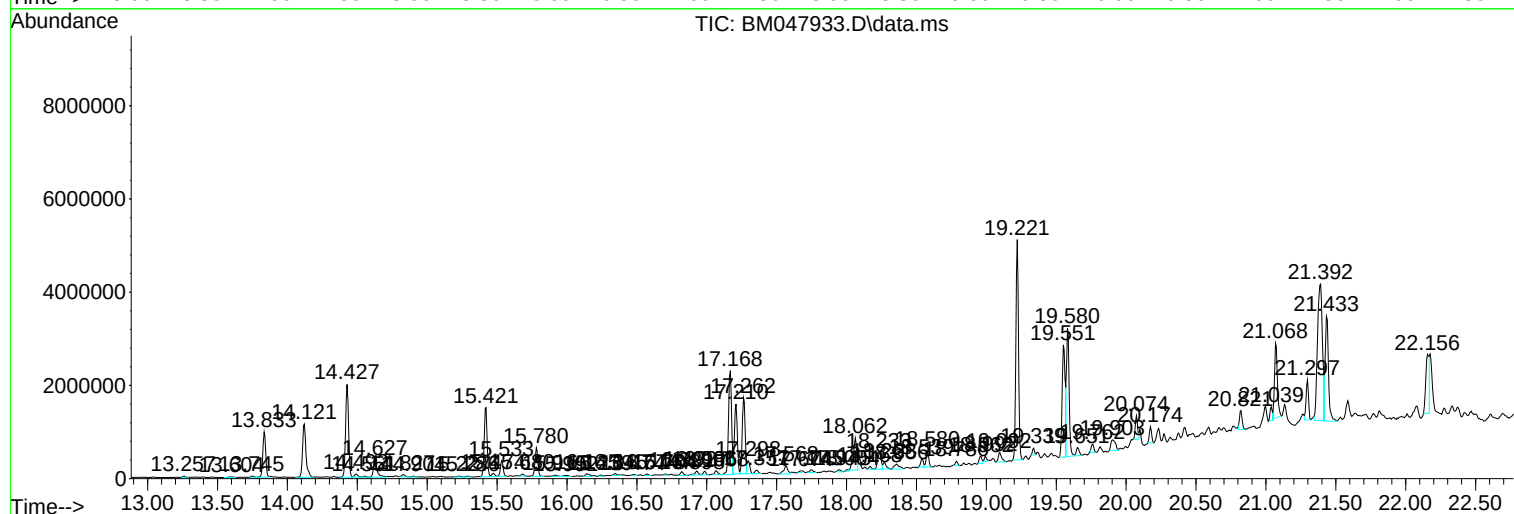
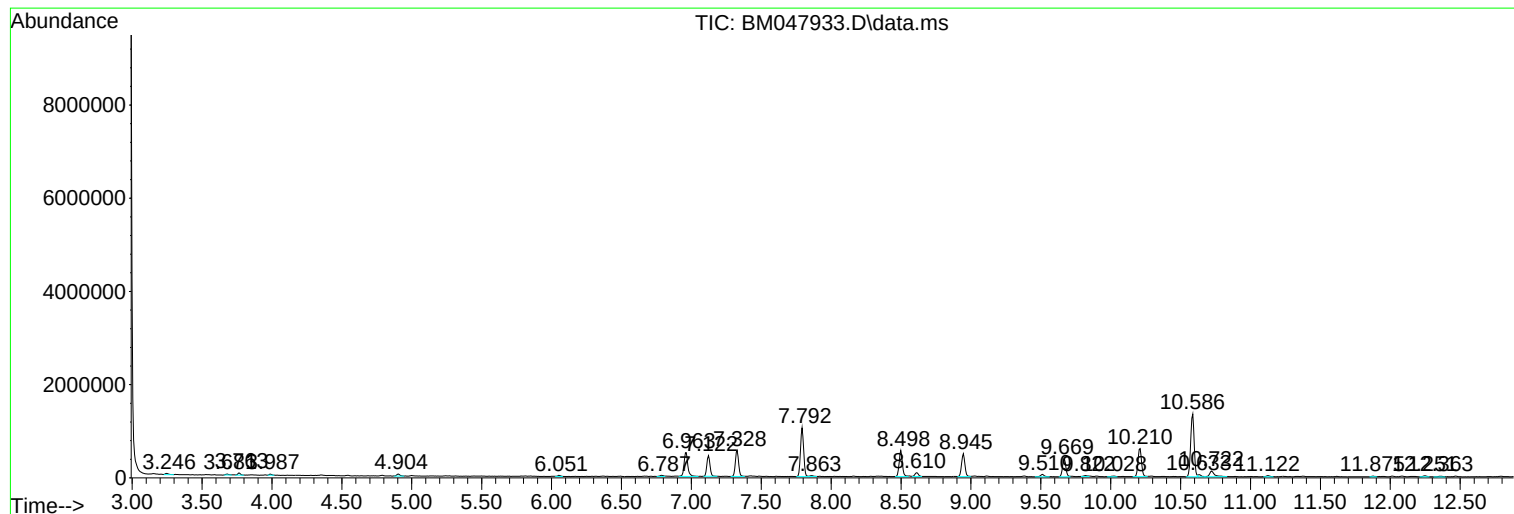
Sum of corrected areas: 84162208

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION

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



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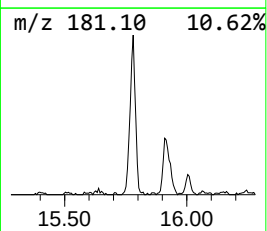
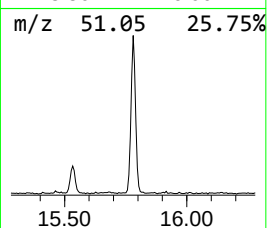
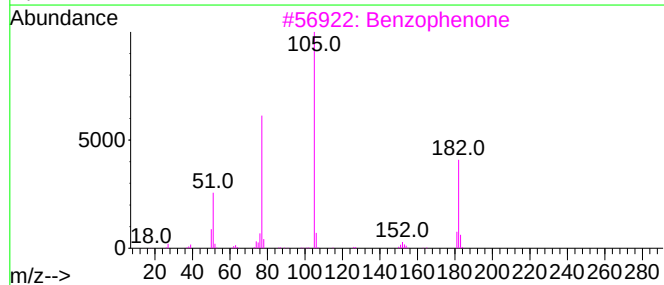
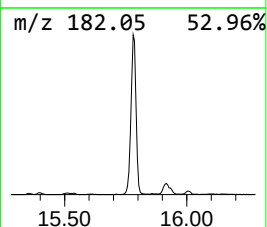
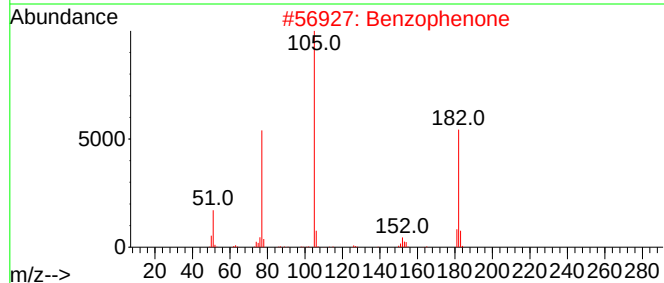
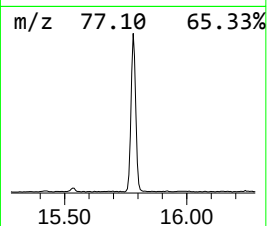
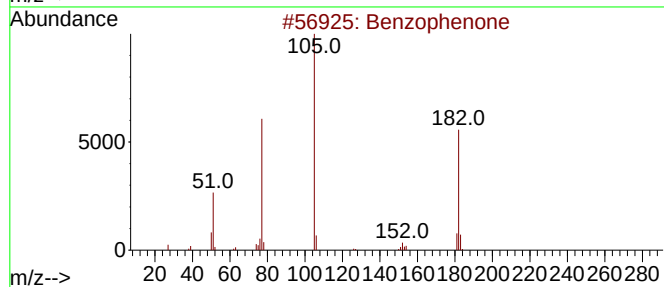
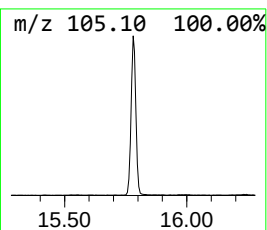
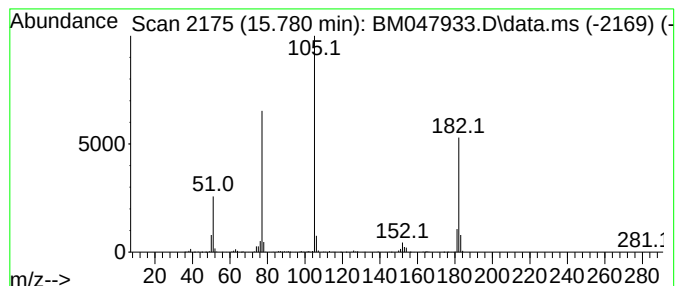
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzophenone Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.780	6.10 ng/ul	873640	Acenaphthene-d10	14.427

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	96
2			Benzophenone	182	C13H10O	000119-61-9	95
3			Benzophenone	182	C13H10O	000119-61-9	94
4			Benzophenone	182	C13H10O	000119-61-9	90
5			Benzoyl isothiocyanate	163	C8H5NOS	000532-55-8	47



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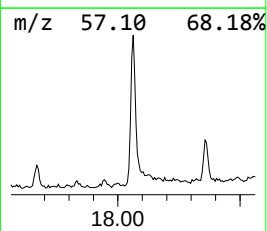
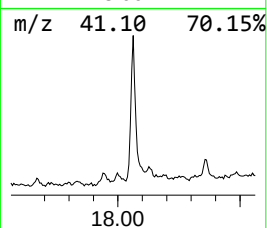
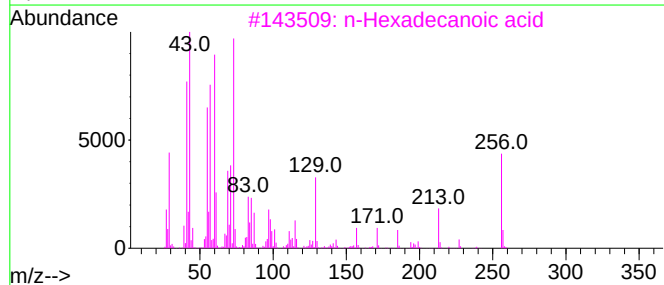
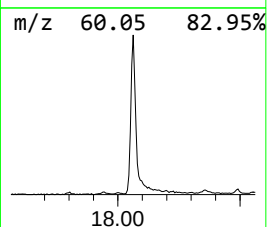
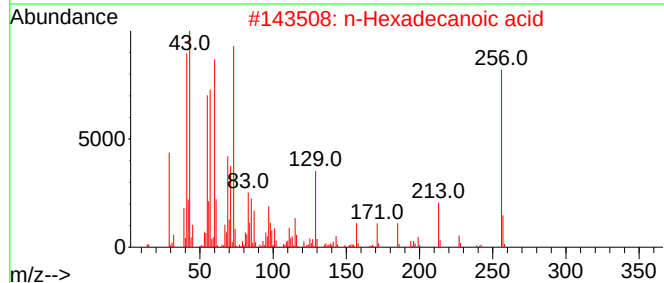
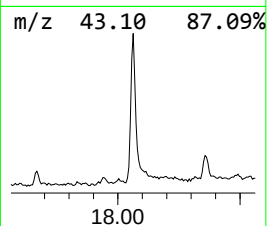
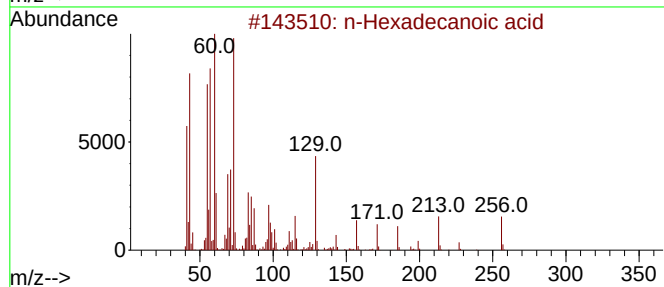
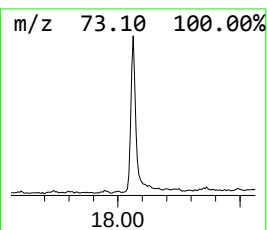
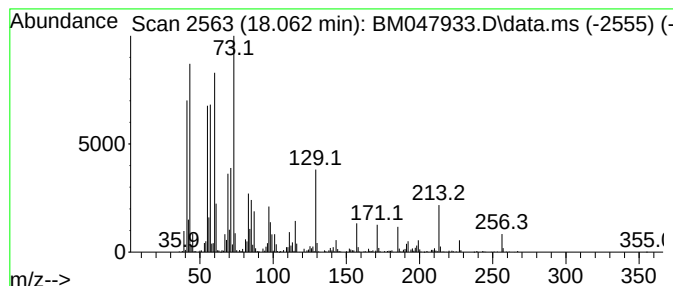
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.062	7.26 ng/ul	1217420	Phenanthrene-d10	17.168

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			Pentadecanoic acid	242	C15H30O2	001002-84-2	93



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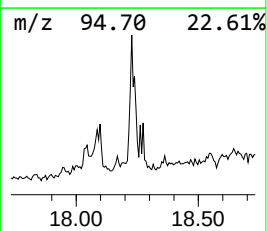
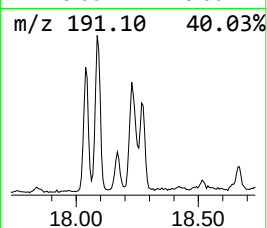
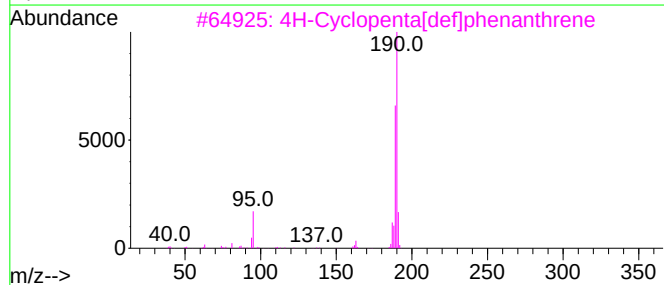
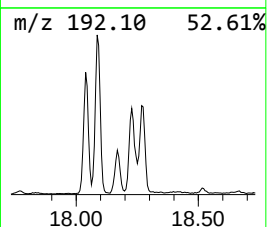
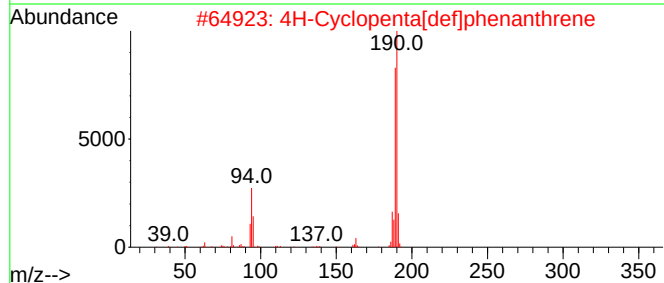
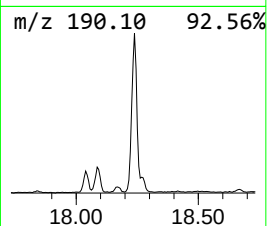
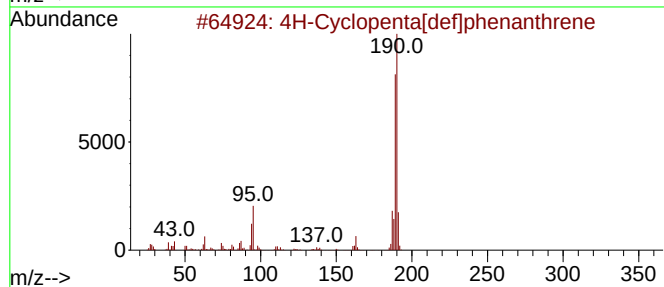
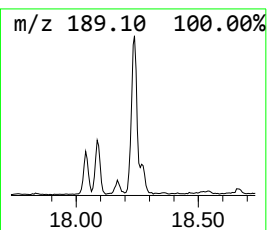
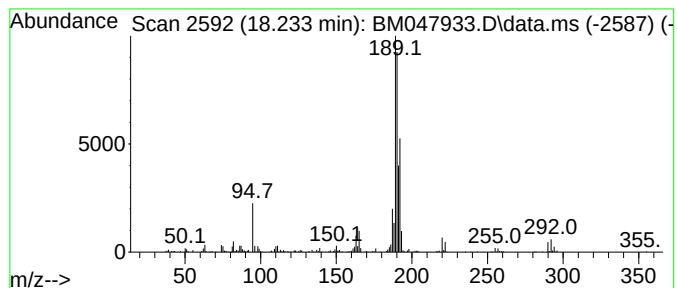
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 4H-Cyclopenta[def]phenanthrene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.233	3.74 ng/ul	627148	Phenanthrene-d10	17.168

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	64
2			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	62
3			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	46
4			Methyl diselenide	190	C2H6Se2	007101-31-7	43
5			Naphtho[1,2-b]norbornadiene	192	C15H12	1000210-14-8	25



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Data File : BM047933.D
Acq On : 09 Oct 2024 04:56
Operator : RC/JU
Sample : P4183-05
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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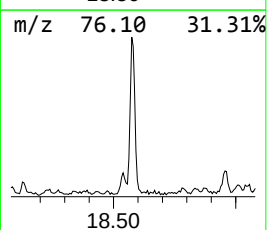
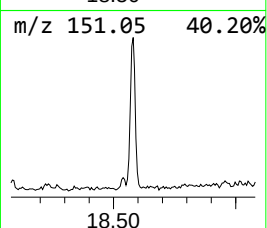
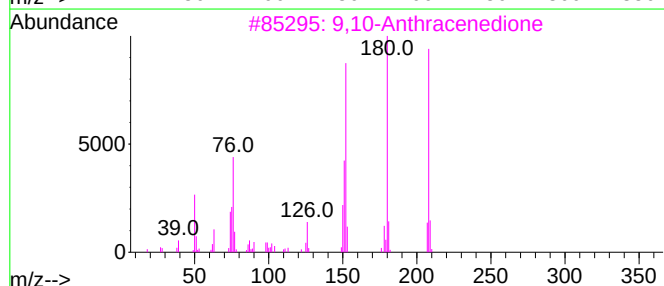
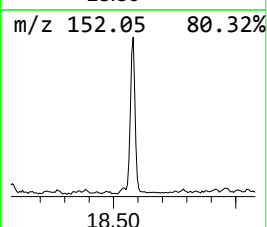
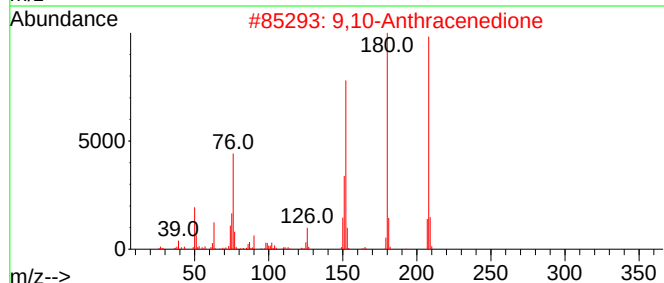
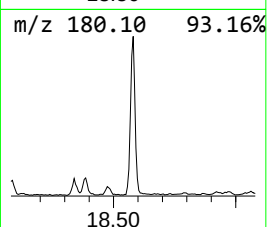
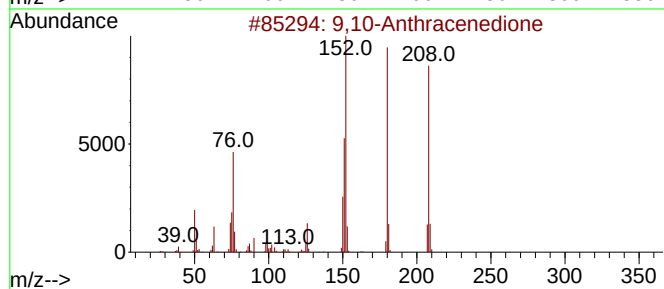
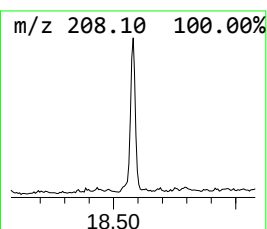
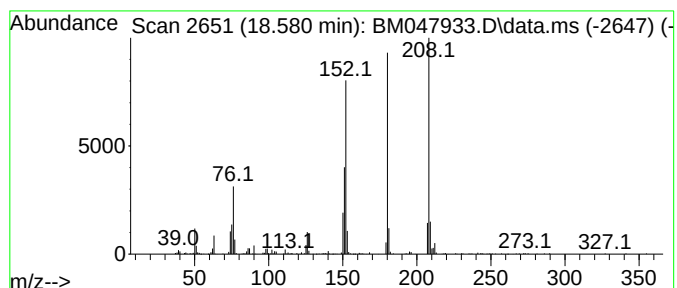
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 9,10-Anthracenedione Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.580	2.74 ng/ul	459165	Phenanthrene-d10	17.168

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9,10-Anthracenedione	208	C14H8O2	000084-65-1	98
2			9,10-Anthracenedione	208	C14H8O2	000084-65-1	95
3			9,10-Anthracenedione	208	C14H8O2	000084-65-1	95
4			9H-Fluoren-9-one	180	C13H8O	000486-25-9	90
5			9,10-Anthracenedione	208	C14H8O2	000084-65-1	89



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100724\
Data File : BM047933.D
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Sample : P4183-05
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Instrument :
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ClientSampleId :
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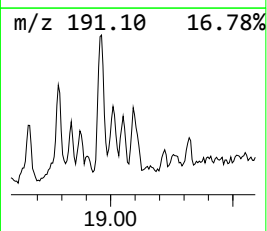
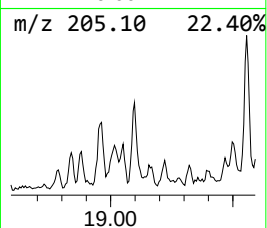
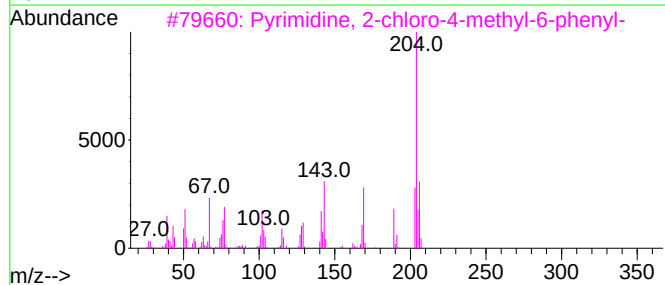
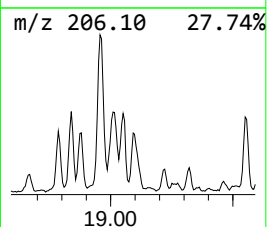
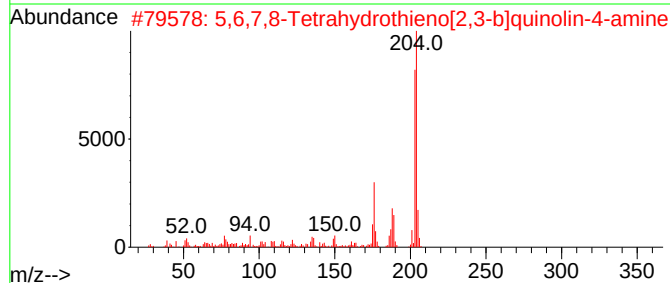
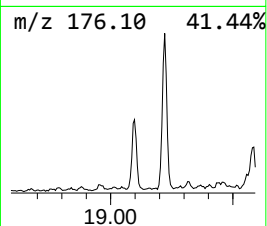
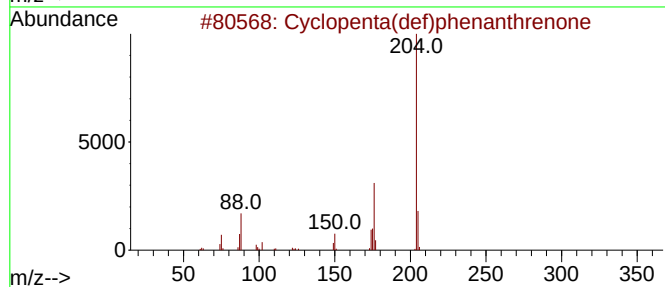
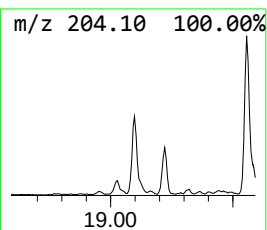
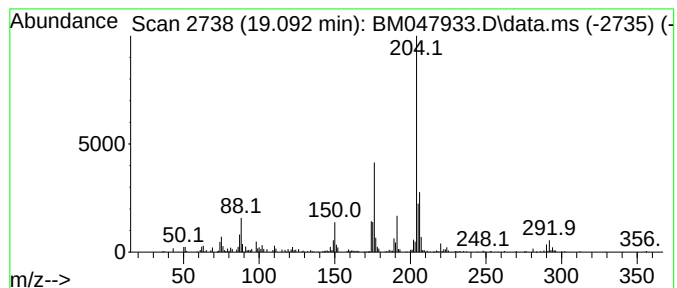
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Cyclopenta(def)phenanthrene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.092	2.01 ng/ul	336737	Phenanthrene-d10	17.168

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopenta(def)phenanthrene	204	C15H8O	005737-13-3	90
2			5,6,7,8-Tetrahydrothieno[2,3-b]q...	204	C11H12N2S	122914-50-5	50
3			Pyrimidine, 2-chloro-4-methyl-6-...	204	C11H9ClN2	032785-40-3	47
4			[1,1'-Biphenyl]-4,4'-dicarbonitrile	204	C14H8N2	001591-30-6	46
5			1,2,4,8-Tetramethylbicyclo[6.3.0...	204	C15H24	137235-51-9	45



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100724\
Data File : BM047933.D
Acq On : 09 Oct 2024 04:56
Operator : RC/JU
Sample : P4183-05
Misc :
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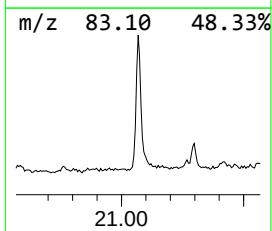
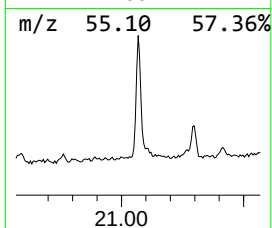
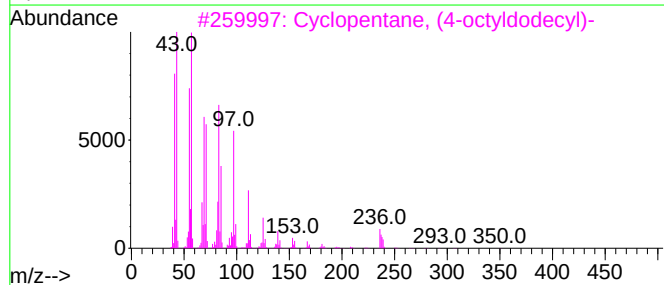
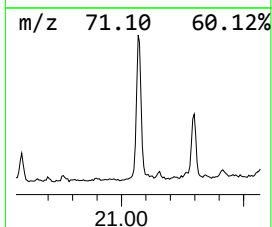
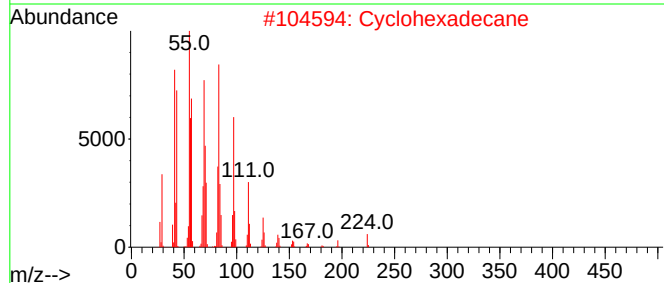
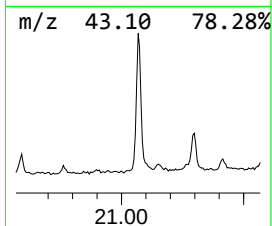
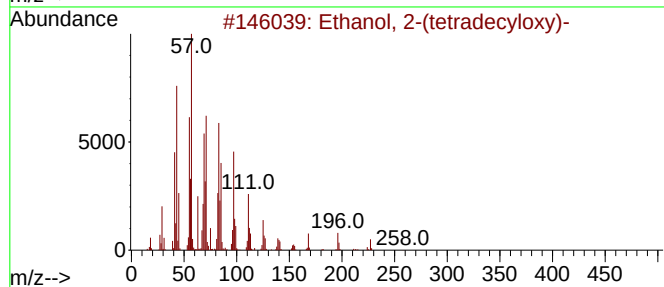
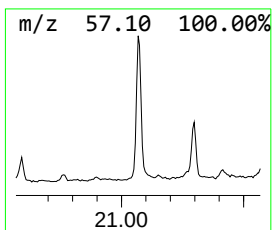
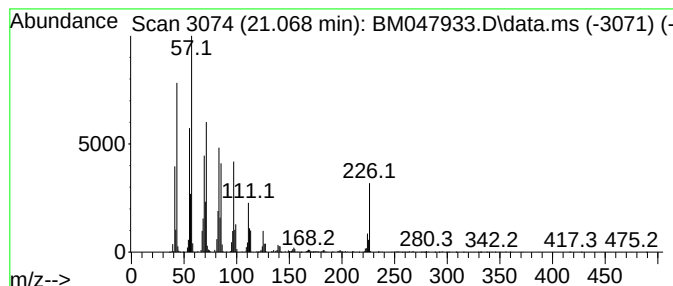
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Ethanol, 2-(tetradecyloxy)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.068	6.25 ng/ul	2290550	Chrysene-d12	21.392	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	99
2	Cyclohexadecane	224	C16H32	000295-65-8	95
3	Cyclopentane, (4-octyldodecyl)-	350	C25H50	005638-09-5	92
4	Oxalic acid, isobutyl hexadecyl ...	370	C22H42O4	1000309-38-1	87
5	Trihexadecyl borate	735	C48H99BO3	002665-11-4	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100724\
Data File : BM047933.D
Acq On : 09 Oct 2024 04:56
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Sample : P4183-05
Misc :
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Instrument :
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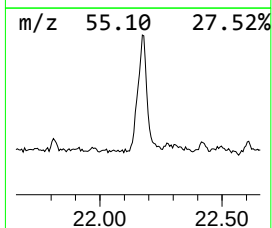
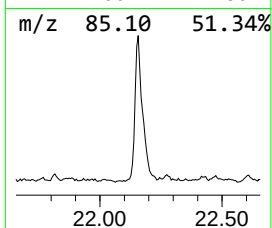
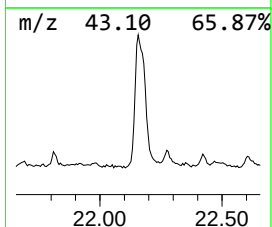
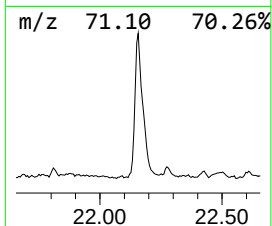
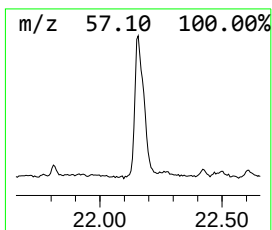
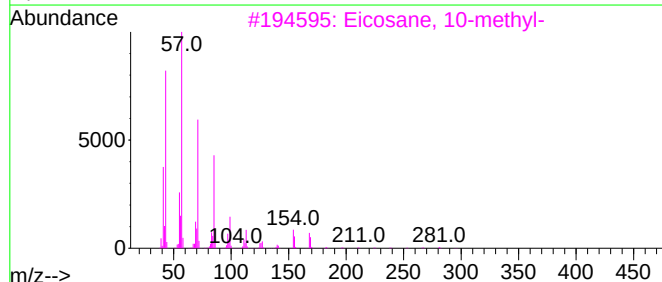
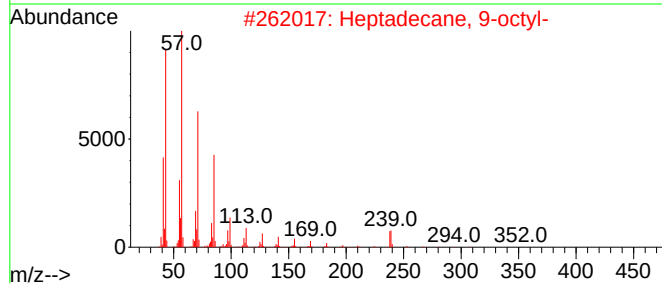
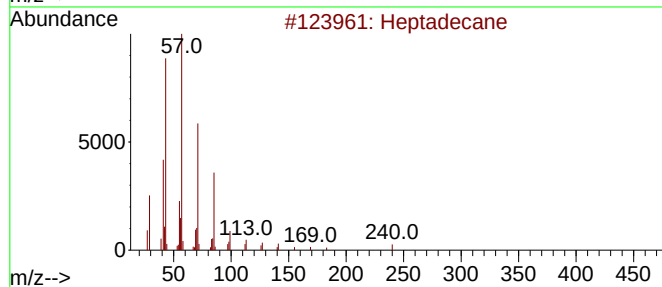
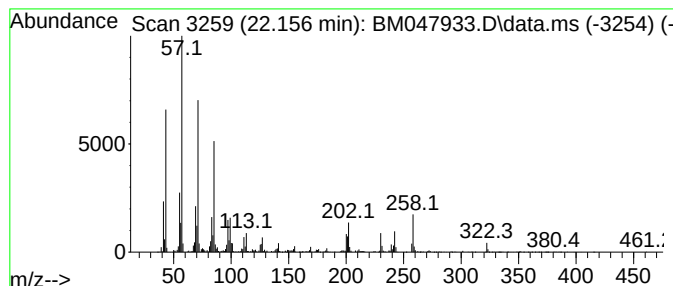
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.156	6.02 ng/ul	2205040	Chrysene-d12	21.392

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecane	240	C17H36	000629-78-7	93
2		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	93
3		Eicosane, 10-methyl-	296	C21H44	054833-23-7	86
4		Heptadecane	240	C17H36	000629-78-7	83
5		Octacosane	394	C28H58	000630-02-4	76



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100724\
Data File : BM047933.D
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Sample : P4183-05
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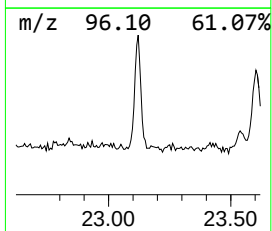
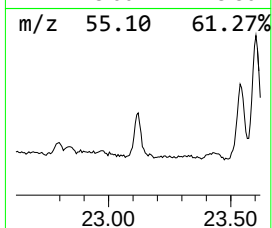
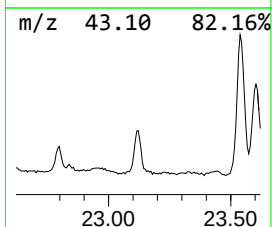
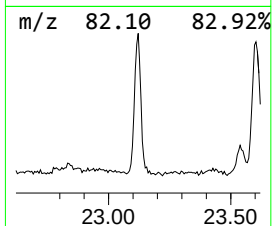
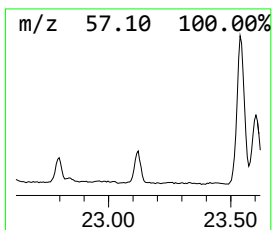
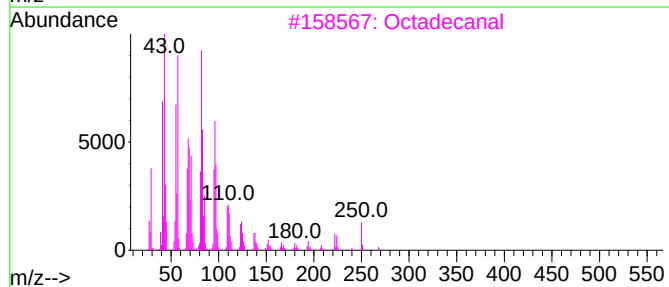
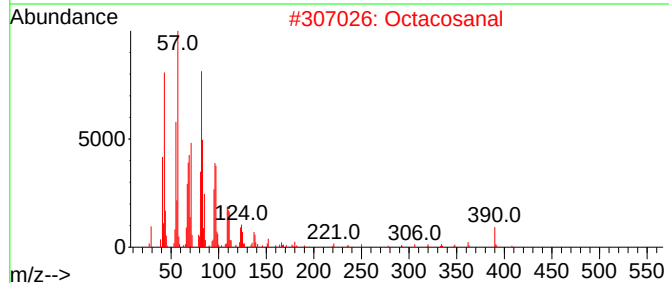
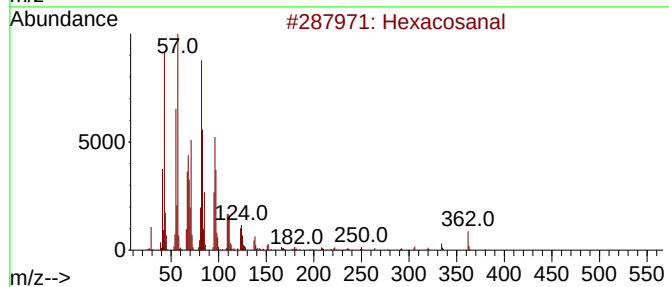
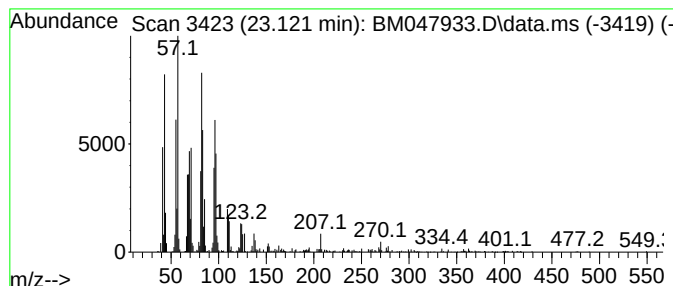
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Hexacosanal Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.121	8.62 ng/ul	1567390	Perylene-d12	24.391

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexacosanal	380	C26H52O	026627-85-0	95
2			Octacosanal	408	C28H56O	022725-64-0	95
3			Octadecanal	268	C18H36O	000638-66-4	93
4			Oxirane, heptadecyl-	282	C19H38O	067860-04-2	93
5			Nonacosanal	422	C29H58O	072934-04-4	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100724\
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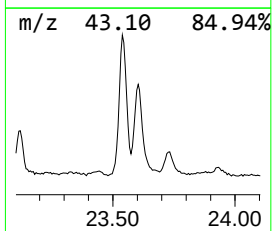
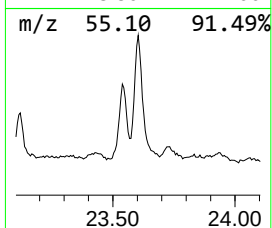
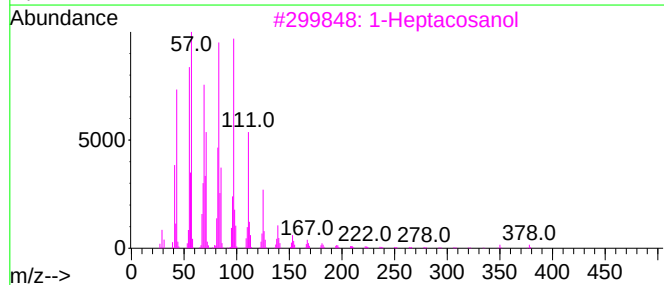
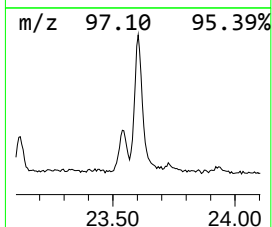
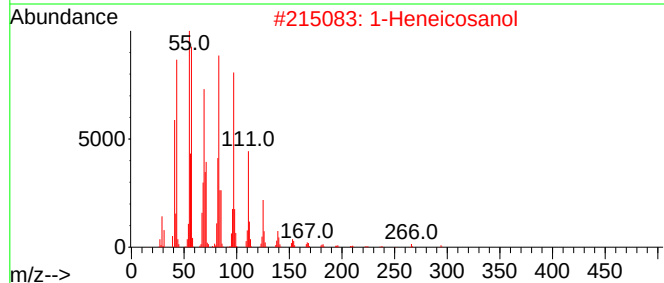
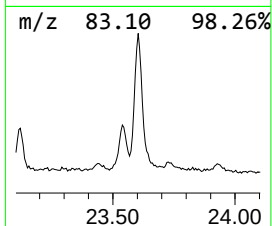
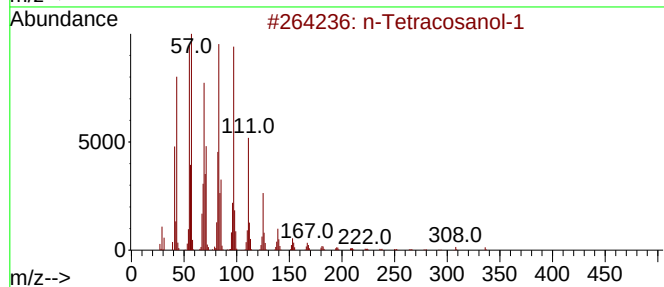
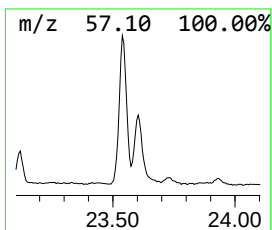
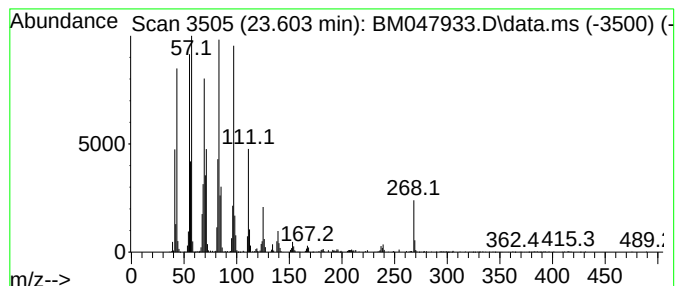
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 n-Tetracosanol-1 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.603	15.52 ng/ul	2820930	Perylene-d12	24.391

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Tetracosanol-1			354	C24H50O	000506-51-4	95
2	1-Heneicosanol			312	C21H44O	015594-90-8	94
3	1-Heptacosanol			396	C27H56O	002004-39-9	94
4	Behenic alcohol			326	C22H46O	000661-19-8	94
5	1-Hexacosene			364	C26H52	018835-33-1	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100724\
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ALS Vial : 23 Sample Multiplier: 1

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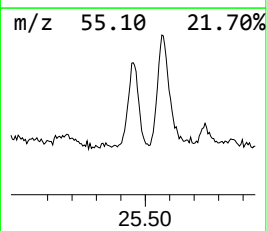
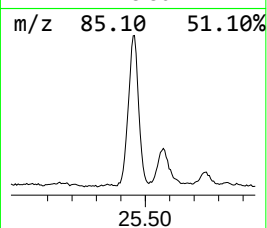
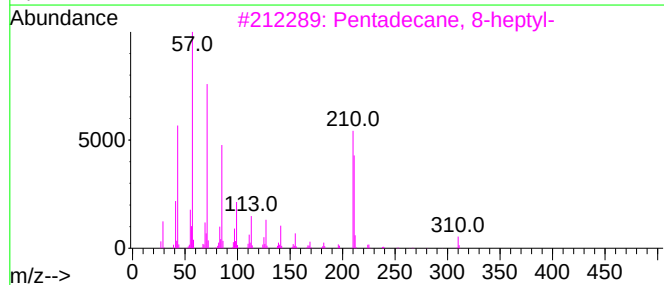
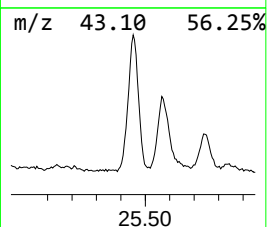
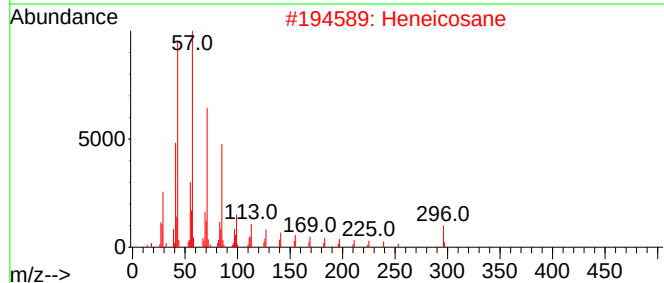
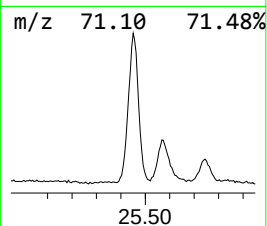
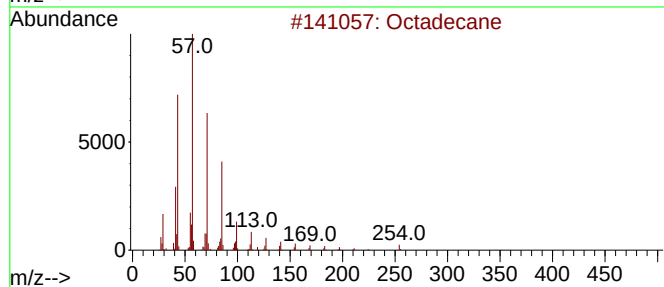
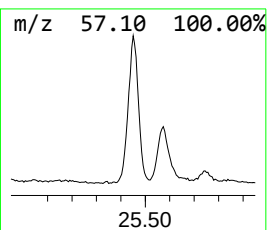
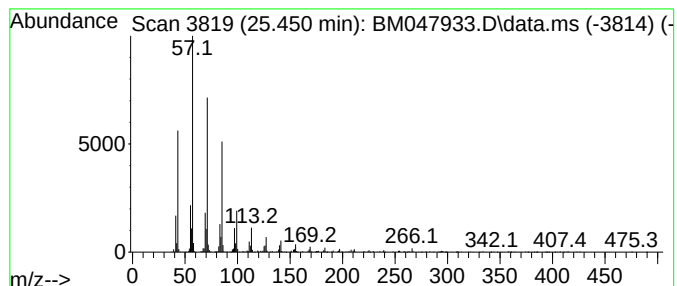
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.450	12.92 ng/ul	2347500	Perylene-d12	24.391

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecane	254	C18H38	000593-45-3	95
2			Heneicosane	296	C21H44	000629-94-7	93
3			Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	91
4			Heneicosane	296	C21H44	000629-94-7	91
5			Dotriacontane, 1-iodo-	576	C32H65I	1000406-32-4	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100724\
Data File : BM047933.D
Acq On : 09 Oct 2024 04:56
Operator : RC/JU
Sample : P4183-05
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
A4EM1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.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
Benzophenone	15.780	6.1	ng/ul	873640	3	14.427	2866090	20.0
n-Hexadecanoic ...	18.062	7.3	ng/ul	1217420	4	17.168	3351920	20.0
4H-Cyclopenta[d...]	18.233	3.7	ng/ul	627148	4	17.168	3351920	20.0
9,10-Anthracene...	18.580	2.7	ng/ul	459165	4	17.168	3351920	20.0
Cyclopenta(def)...	19.092	2.0	ng/ul	336737	4	17.168	3351920	20.0
Ethanol, 2-(tet...	21.068	6.3	ng/ul	2290550	5	21.392	7325410	20.0
(DEL) Alkane: S...	22.156	6.0	ng/ul	2205040	5	21.392	7325410	20.0
Hexacosanal	23.121	8.6	ng/ul	1567390	6	24.391	3634750	20.0
n-Tetracosanol-1	23.603	15.5	ng/ul	2820930	6	24.391	3634750	20.0
(DEL) Alkane: S...	25.450	12.9	ng/ul	2347500	6	24.391	3634750	20.0