

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101424\
 Data File : BM048027.D
 Acq On : 14 Oct 2024 14:15
 Operator : RC/JU
 Sample : P4239-06
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4EG2

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: BM048027.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.240	40	43	50	rVB	24523	32535	0.44%	0.049%
2	3.663	110	115	122	rBV2	86323	131212	1.79%	0.196%
3	3.757	126	131	139	rVV	398628	512375	6.97%	0.765%
4	3.834	140	144	154	rVB	43904	67077	0.91%	0.100%
5	3.981	165	169	174	rVB	21929	30284	0.41%	0.045%
6	4.534	259	263	269	rVB3	67987	92393	1.26%	0.138%
7	4.781	300	305	315	rVB3	29651	51110	0.70%	0.076%
8	4.893	320	324	332	rVB2	58495	90308	1.23%	0.135%
9	4.987	335	340	353	rBV	59315	104999	1.43%	0.157%
10	6.045	514	520	527	rVB2	39829	64838	0.88%	0.097%
11	6.957	666	675	688	rBV	322157	601718	8.19%	0.898%
12	7.116	696	702	713	rBV	281180	461270	6.28%	0.689%
13	7.316	730	736	746	rBV	368049	595962	8.11%	0.890%
14	7.781	808	815	824	rBV	1100699	1812257	24.66%	2.705%
15	8.151	871	878	884	rVB	11885	24662	0.34%	0.037%
16	8.492	930	936	943	rBV	231649	404895	5.51%	0.604%
17	8.604	950	955	963	rVB2	56053	101418	1.38%	0.151%
18	8.939	1005	1012	1022	rVV	303884	544446	7.41%	0.813%
19	9.175	1045	1052	1059	rBV	35745	56993	0.78%	0.085%
20	9.369	1079	1085	1090	rBV5	15482	29933	0.41%	0.045%
21	9.657	1127	1134	1141	rBV	253895	441820	6.01%	0.660%
22	9.822	1158	1162	1169	rBV2	19722	39320	0.53%	0.059%
23	10.204	1219	1227	1236	rBV	352105	662393	9.01%	0.989%
24	10.575	1282	1290	1296	rBV	1489722	2490551	33.88%	3.718%
25	10.622	1296	1298	1306	rVB	45120	69462	0.95%	0.104%
26	10.722	1309	1315	1323	rBV	36773	95306	1.30%	0.142%
27	11.357	1416	1423	1429	rBV5	10921	29049	0.40%	0.043%
28	11.427	1429	1435	1440	rVB	15045	28402	0.39%	0.042%
29	12.239	1569	1573	1582	rBV4	24761	48492	0.66%	0.072%
30	12.363	1587	1594	1603	rBV7	7825	19731	0.27%	0.029%
31	12.457	1603	1610	1615	rVB2	15561	27975	0.38%	0.042%
32	13.598	1797	1804	1809	rBV8	11569	25537	0.35%	0.038%
33	13.745	1823	1829	1833	rBV3	20954	35283	0.48%	0.053%
34	13.827	1833	1843	1849	rBV	671778	988193	13.44%	1.475%
35	14.110	1880	1891	1906	rBV2	806603	1435717	19.53%	2.143%
36	14.263	1910	1917	1925	rBV2	22276	58934	0.80%	0.088%

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Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
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37	14.421	1933	1944	1950	rBV	2110610	3191834	43.43%	4.765%
38	14.486	1950	1955	1961	rVB	91197	145541	1.98%	0.217%
39	14.545	1961	1965	1973	rVB5	24943	43415	0.59%	0.065%
40	14.633	1973	1980	1984	rBV	150049	297743	4.05%	0.444%

41	14.668	1984	1986	1992	rVB2	73272	104703	1.42%	0.156%
42	14.774	2000	2004	2007	rBV4	17965	29560	0.40%	0.044%
43	14.821	2007	2012	2019	rVV	51059	104862	1.43%	0.157%
44	14.886	2019	2023	2029	rVB4	40656	64869	0.88%	0.097%
45	15.027	2043	2047	2053	rBV8	10123	22772	0.31%	0.034%

46	15.210	2072	2078	2086	rBV9	15660	41893	0.57%	0.063%
47	15.409	2106	2112	2118	rBV	1078794	1552549	21.12%	2.318%
48	15.468	2118	2122	2128	rVV	96688	142569	1.94%	0.213%
49	15.533	2128	2133	2140	rVV	226010	338906	4.61%	0.506%
50	15.592	2140	2143	2146	rVV4	18028	23819	0.32%	0.036%

51	15.627	2146	2149	2152	rVV4	14262	20716	0.28%	0.031%
52	15.674	2152	2157	2162	rVB	50963	74458	1.01%	0.111%
53	15.774	2168	2174	2183	rVB2	269988	418119	5.69%	0.624%
54	15.904	2193	2196	2204	rVB	30856	62794	0.85%	0.094%
55	15.986	2204	2210	2217	rBV8	24978	49332	0.67%	0.074%

56	16.139	2231	2236	2242	rBV5	44307	78816	1.07%	0.118%
57	16.274	2256	2259	2265	rBV8	13040	25784	0.35%	0.038%
58	16.392	2275	2279	2283	rVB2	38925	51468	0.70%	0.077%
59	16.568	2305	2309	2312	rBV6	16718	21176	0.29%	0.032%
60	16.698	2327	2331	2334	rBV4	24347	33429	0.45%	0.050%

61	16.739	2335	2338	2342	rVB5	19363	24683	0.34%	0.037%
62	16.815	2347	2351	2358	rVB	79020	113752	1.55%	0.170%
63	16.915	2365	2368	2374	rVV3	54817	87740	1.19%	0.131%
64	16.980	2375	2379	2383	rVB	78529	104140	1.42%	0.155%
65	17.162	2400	2410	2413	rBV	2449927	3689893	50.20%	5.508%

66	17.204	2413	2417	2421	rVV	1523963	2220896	30.22%	3.315%
67	17.256	2421	2426	2430	rVV	1153725	1702922	23.17%	2.542%
68	17.292	2430	2432	2436	rVV	307022	347592	4.73%	0.519%
69	17.351	2437	2442	2448	rVB4	51361	105633	1.44%	0.158%
70	17.562	2471	2478	2483	rBV	169611	286070	3.89%	0.427%

71	17.739	2505	2508	2518	rVB3	51055	98390	1.34%	0.147%
72	17.992	2548	2551	2553	rBV3	34762	41202	0.56%	0.062%
73	18.051	2554	2561	2564	rBV2	620831	1093583	14.88%	1.632%
74	18.162	2577	2580	2585	rVB	67626	90084	1.23%	0.134%
75	18.227	2585	2591	2595	rBV3	335532	608277	8.28%	0.908%

76	18.262	2595	2597	2604	rVB	141636	195820	2.66%	0.292%
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77	18.345	2606	2611	2614	rBV5	57377	111807	1.52%	0.167%
78	18.533	2639	2643	2646	rBV	184142	226401	3.08%	0.338%
79	18.574	2646	2650	2656	rVB	286663	388402	5.28%	0.580%
80	18.780	2682	2685	2689	rVB2	70996	81085	1.10%	0.121%
81	18.950	2710	2714	2718	rBV2	150093	249970	3.40%	0.373%
82	19.092	2734	2738	2746	rBV	161456	286057	3.89%	0.427%
83	19.215	2753	2759	2766	rVB	3822565	5103611	69.44%	7.619%
84	19.333	2776	2779	2781	rVB	131089	123726	1.68%	0.185%
85	19.550	2810	2816	2818	rBV2	1833590	2987756	40.65%	4.460%
86	19.574	2818	2820	2828	rVV	2699125	3394335	46.18%	5.067%
87	19.650	2830	2833	2839	rVB2	150491	223720	3.04%	0.334%
88	19.756	2848	2851	2856	rVB	138994	178300	2.43%	0.266%
89	19.903	2871	2876	2882	rBV3	180378	448175	6.10%	0.669%
90	20.068	2901	2904	2911	rVB3	439805	711405	9.68%	1.062%
91	20.168	2918	2921	2925	rBV	269891	314484	4.28%	0.469%
92	20.815	3028	3031	3037	rVB	367824	517680	7.04%	0.773%
93	21.033	3065	3068	3070	rBV	249695	293509	3.99%	0.438%
94	21.062	3070	3073	3081	rVB4	857463	1383760	18.83%	2.066%
95	21.291	3108	3112	3116	rVB	736534	896187	12.19%	1.338%
96	21.386	3121	3128	3133	rVV3	3134988	7350148	100.00%	10.972%
97	21.433	3133	3136	3148	rVB	1779733	2682420	36.49%	4.004%
98	22.144	3253	3257	3259	rBV3	534219	825851	11.24%	1.233%
99	23.438	3471	3477	3484	rBV	1368459	3743277	50.93%	5.588%
100	24.385	3631	3638	3649	rVB	1729801	4408557	59.98%	6.581%

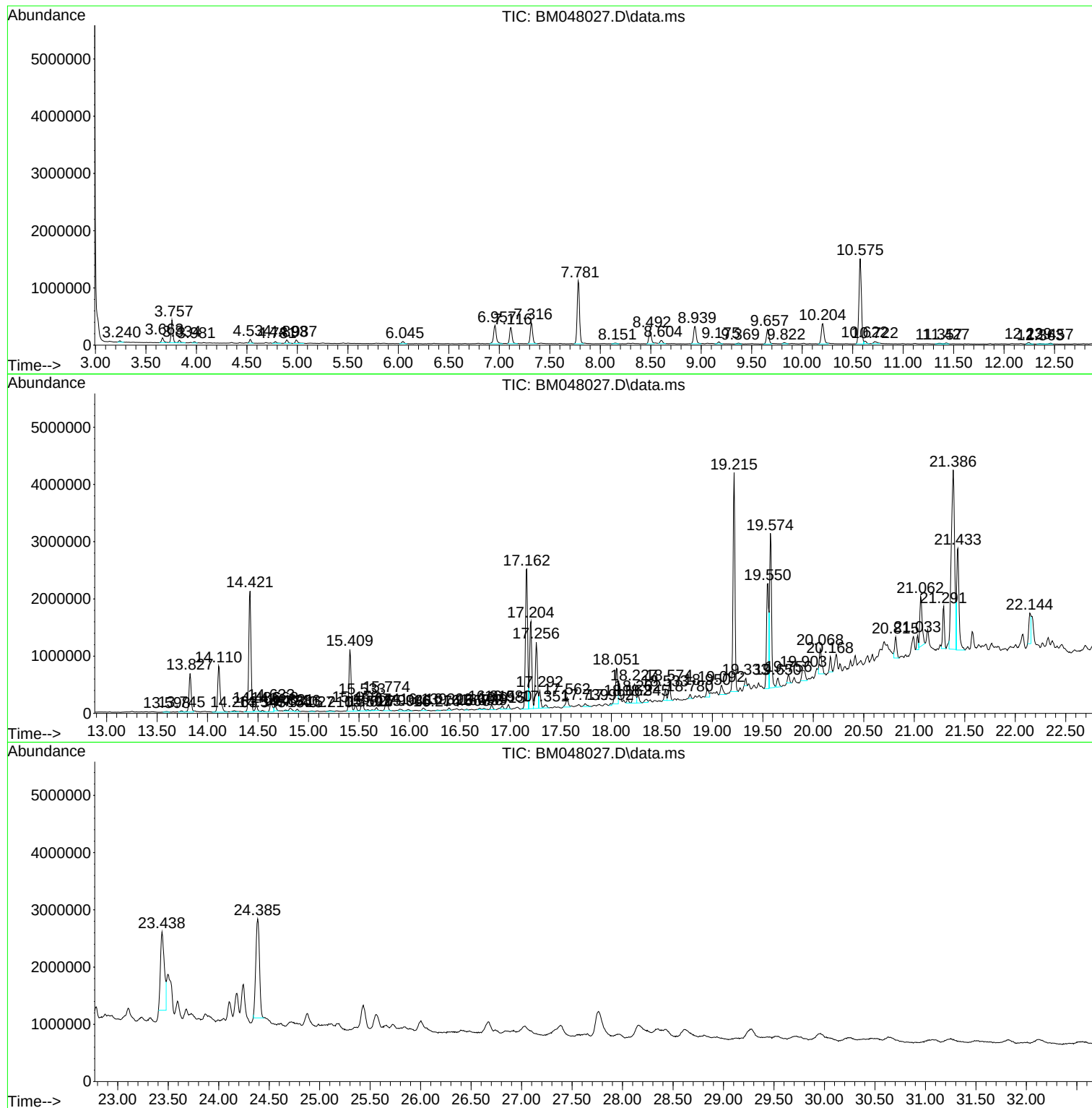
Sum of corrected areas: 66989277

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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



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101424\
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Instrument :
BNA_M
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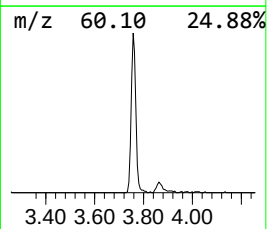
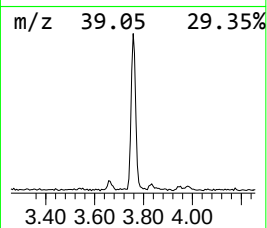
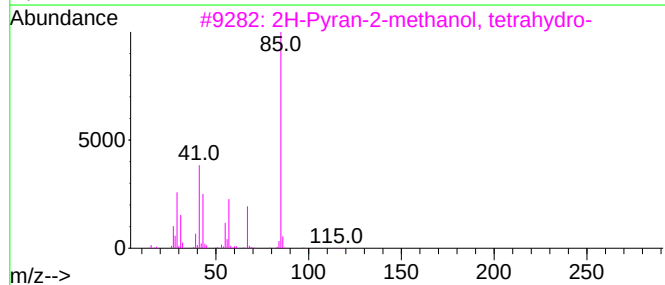
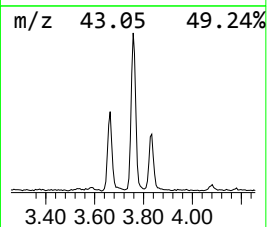
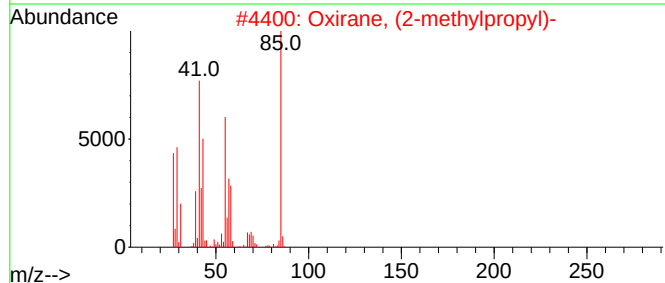
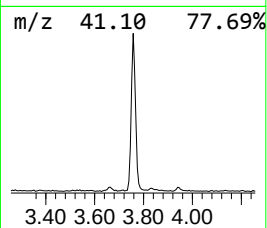
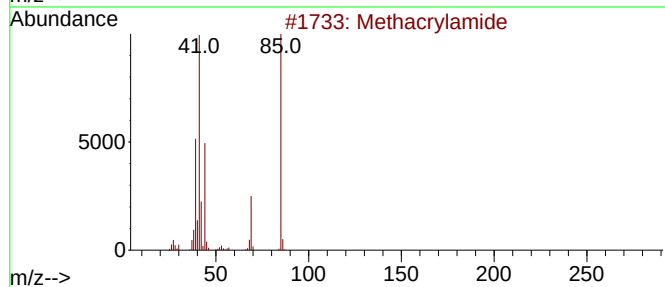
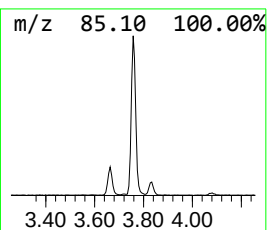
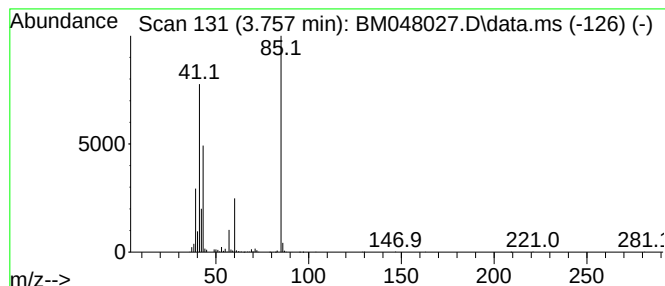
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 unknown-01 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.757	5.65 ng/ul	512375	1,4-Dichlorobenzene-d4	7.786

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methacrylamide	85	C4H7NO	000079-39-0	36
2			Oxirane, (2-methylpropyl)-	100	C6H12O	023850-78-4	36
3			2H-Pyran-2-methanol, tetrahydro-	116	C6H12O2	000100-72-1	9
4			Azetidine, 1-nitroso-	86	C3H6N2O	015216-10-1	9
5			3-Butenoic acid	86	C4H6O2	000625-38-7	9



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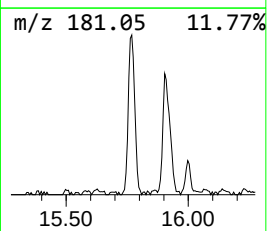
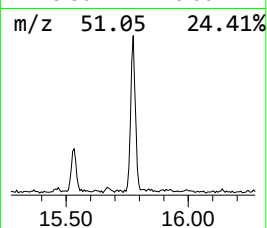
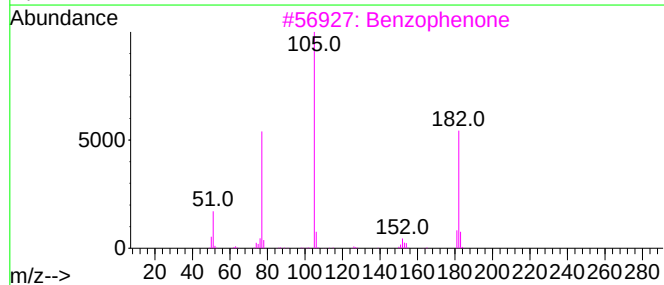
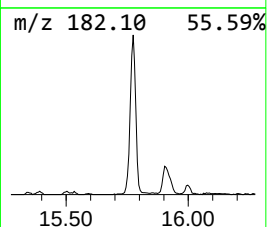
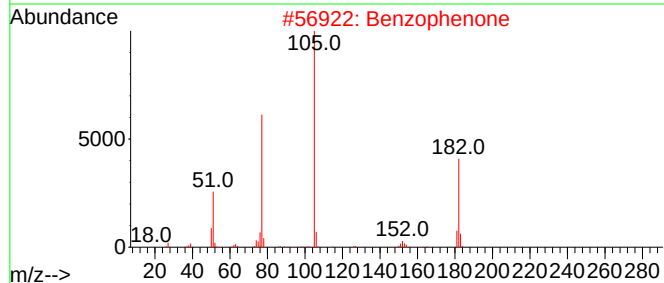
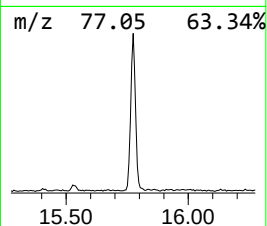
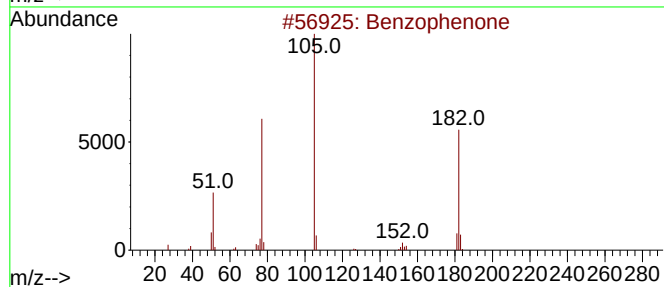
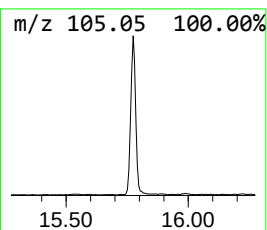
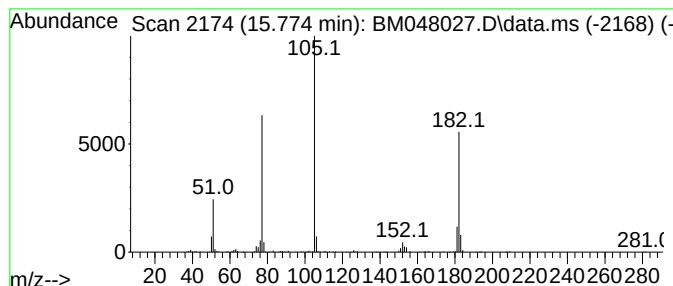
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzophenone Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.774	2.62 ng/ul	418119	Acenaphthene-d10	14.421

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	95
4			Benzophenone	182	C13H10O	000119-61-9	90
5			1-Propanone, 2-bromo-1-phenyl-	212	C9H9BrO	002114-00-3	47



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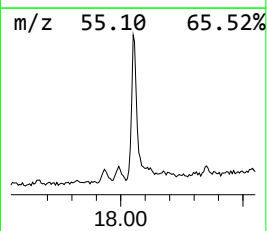
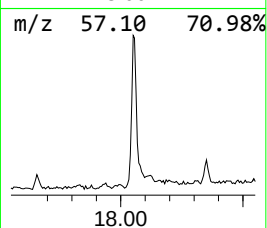
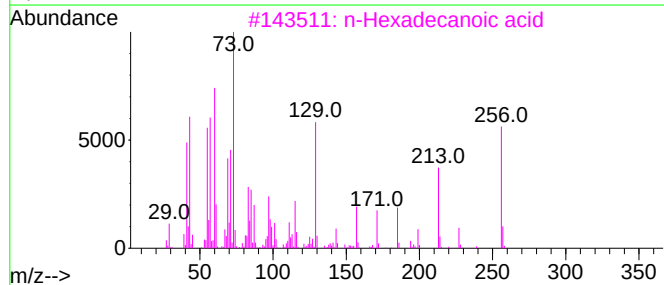
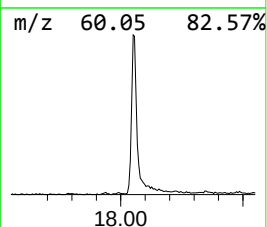
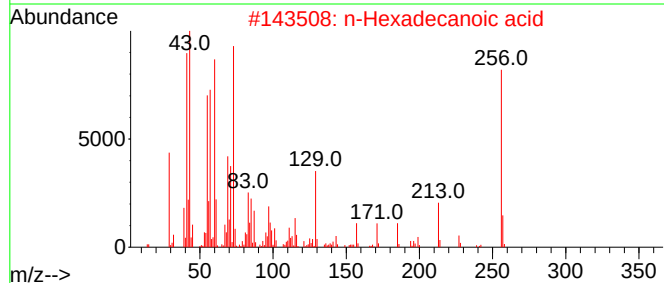
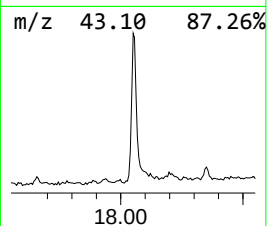
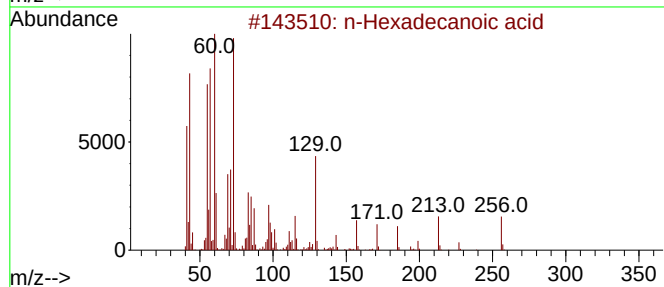
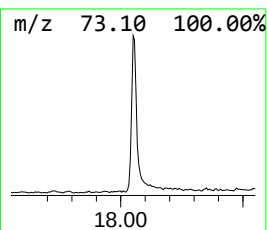
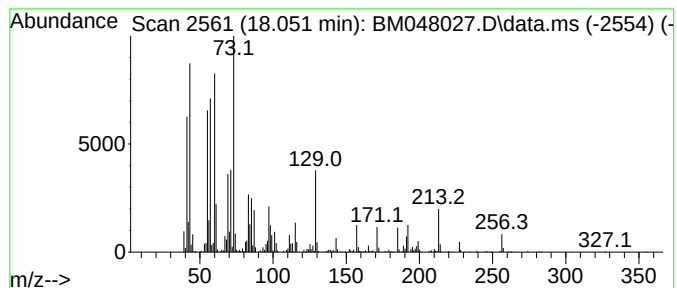
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 n-Hexadecanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.051	5.93 ng/ul	1093580	Phenanthrene-d10	17.162

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	98
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
5			Tridecanoic acid	214	C13H26O2	000638-53-9	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101424\
Data File : BM048027.D
Acq On : 14 Oct 2024 14:15
Operator : RC/JU
Sample : P4239-06
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
A4EG2

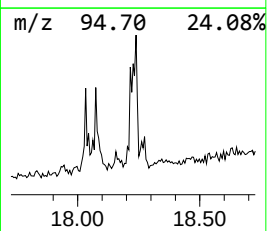
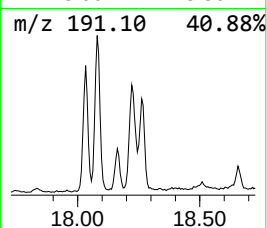
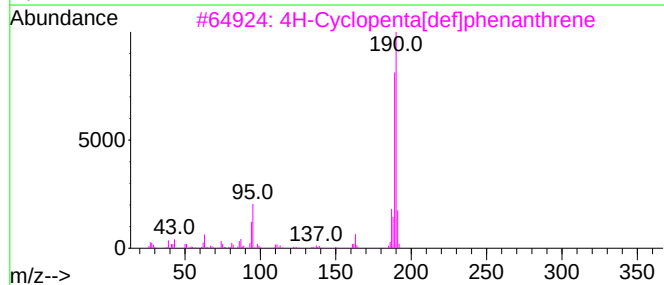
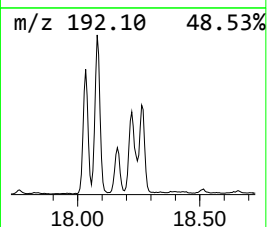
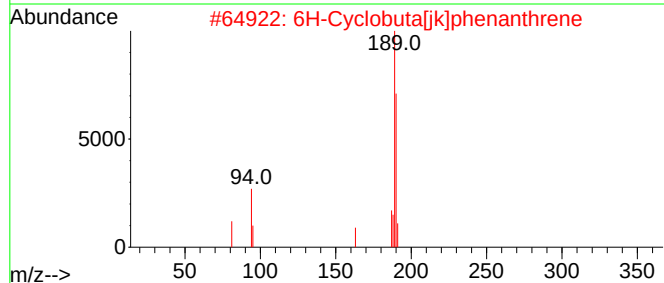
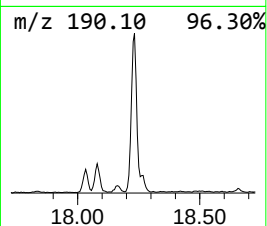
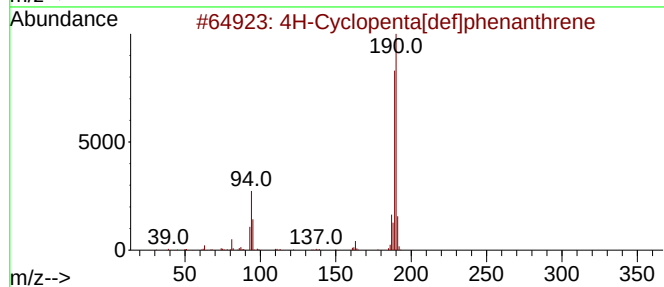
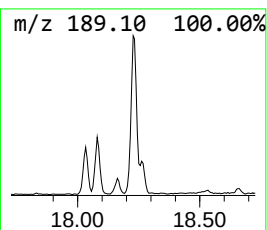
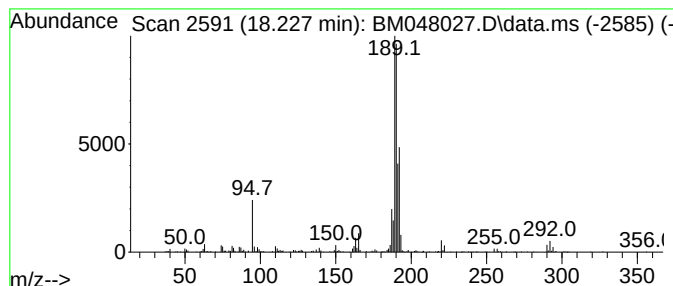
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.227	3.30 ng/ul	608277	Phenanthrene-d10	17.162

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	70
2			6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	53
3			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	50
4			Carbonic acid, ethyl 2-formyl-4,...	262	C10H8Cl2O4	1000331-34-4	50
5			Pyrazole-4-carboxaldehyde, 3-(4-...	190	C10H7FN2O	306936-57-2	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101424\
Data File : BM048027.D
Acq On : 14 Oct 2024 14:15
Operator : RC/JU
Sample : P4239-06
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
A4EG2

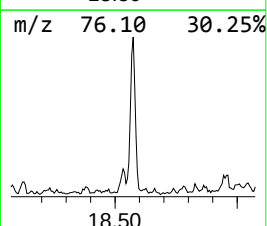
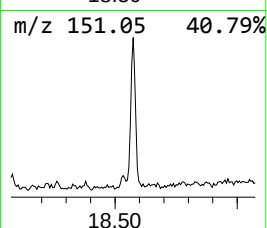
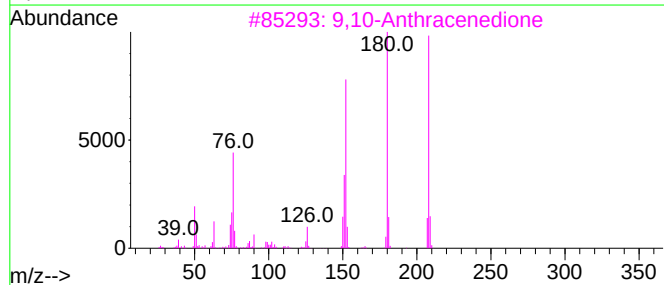
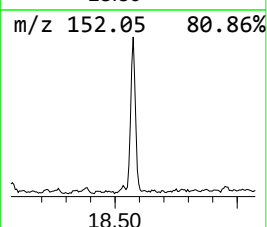
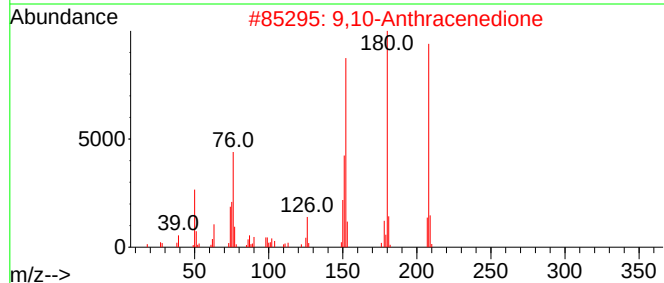
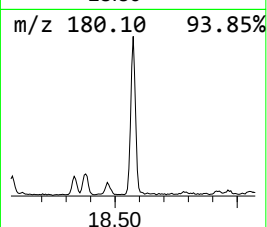
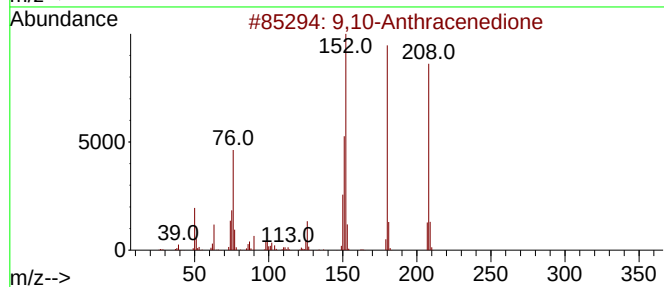
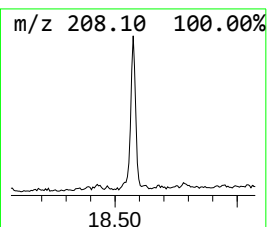
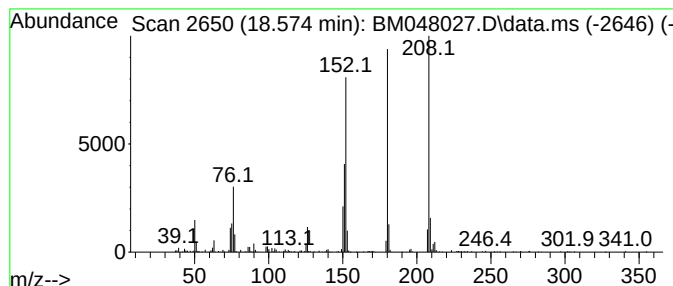
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 9,10-Anthracenedione Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.574	2.11 ng/ul	388402	Phenanthrene-d10	17.162

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	94
3	9,10-Anthracenedione			208	C14H8O2	000084-65-1	93
4	9,10-Anthracenedione			208	C14H8O2	000084-65-1	93
5	9H-Fluoren-9-one			180	C13H8O	000486-25-9	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101424\
Data File : BM048027.D
Acq On : 14 Oct 2024 14:15
Operator : RC/JU
Sample : P4239-06
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
A4EG2

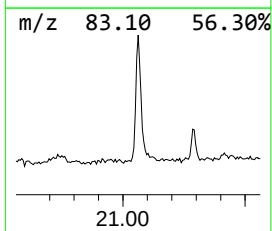
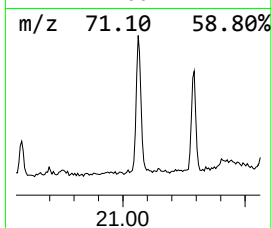
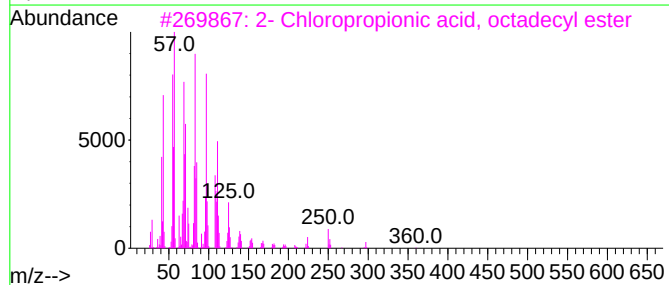
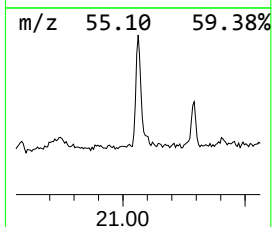
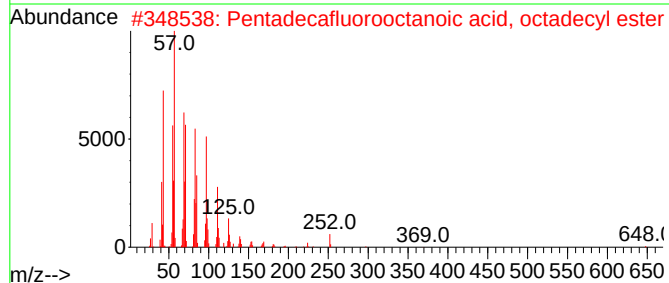
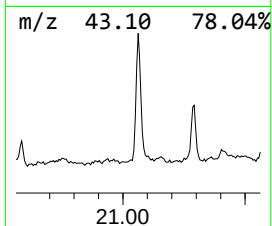
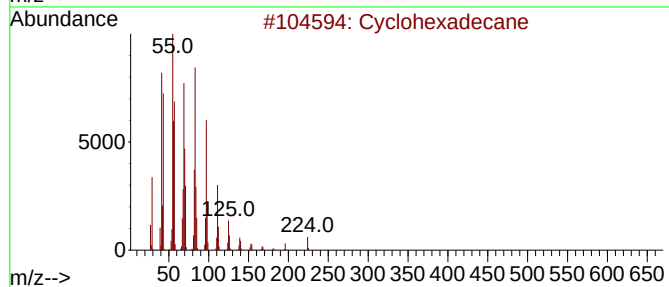
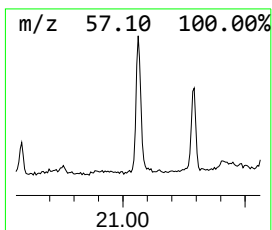
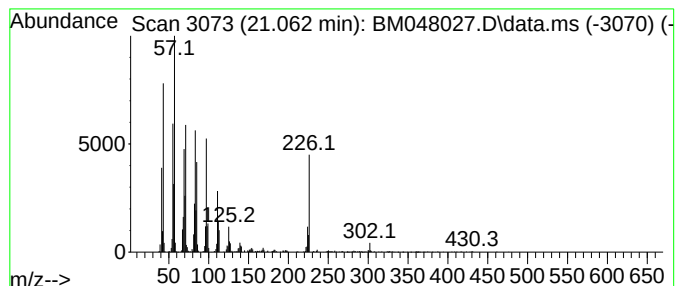
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Cyclic21.062 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.062	3.77 ng/ul	1383760	Chrysene-d12	21.386

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexadecane	224	C16H32	000295-65-8	97
2			Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	93
3			2- Chloropropionic acid, octadec...	360	C21H41ClO2	088104-31-8	92
4			Trihexadecyl borate	735	C48H99B03	002665-11-4	81
5			Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101424\
Data File : BM048027.D
Acq On : 14 Oct 2024 14:15
Operator : RC/JU
Sample : P4239-06
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
A4EG2

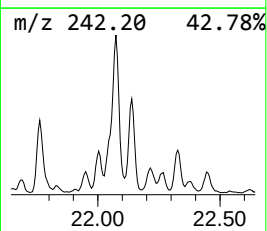
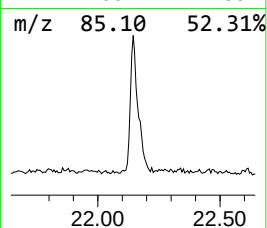
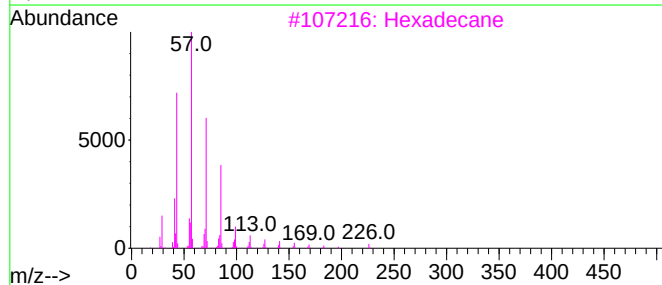
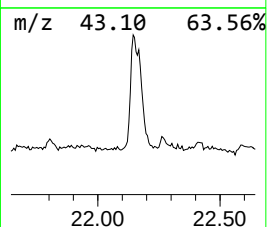
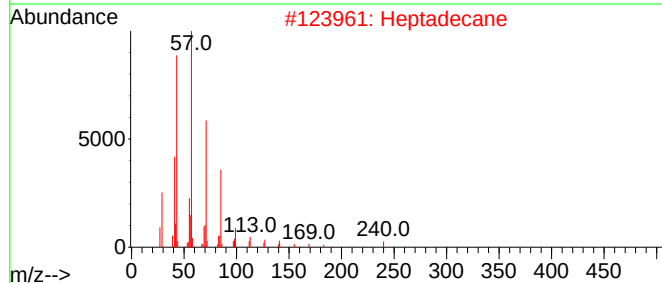
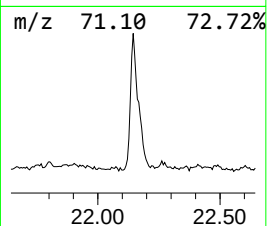
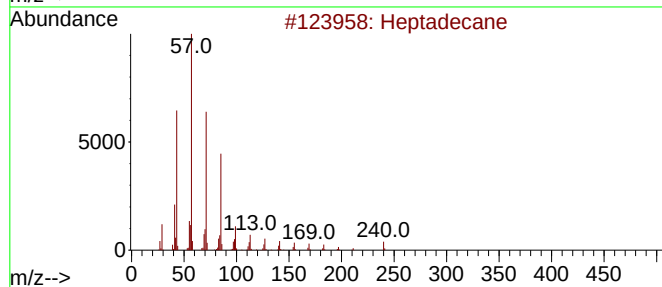
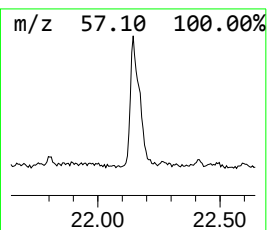
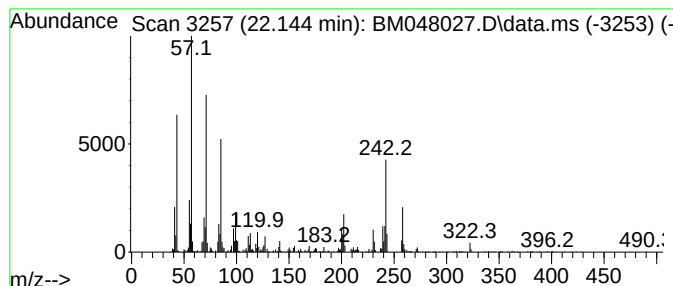
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.144	2.25 ng/ul	825851	Chrysene-d12	21.386

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecane	240	C17H36	000629-78-7	95
2		Heptadecane	240	C17H36	000629-78-7	83
3		Hexadecane	226	C16H34	000544-76-3	83
4		Tetracosane	338	C24H50	000646-31-1	42
5		Octadecane	254	C18H38	000593-45-3	42



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101424\
Data File : BM048027.D
Acq On : 14 Oct 2024 14:15
Operator : RC/JU
Sample : P4239-06
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
A4EG2

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
unknown-01	3.757	5.7	ng/u1	512375	1	7.786	1812260	20.0
Benzophenone	15.774	2.6	ng/u1	418119	3	14.421	3191830	20.0
n-Hexadecanoic ...	18.051	5.9	ng/u1	1093580	4	17.162	3689890	20.0
4H-Cyclopenta[d...]	18.227	3.3	ng/u1	608277	4	17.162	3689890	20.0
9,10-Anthracene...	18.574	2.1	ng/u1	388402	4	17.162	3689890	20.0
(DEL) Alkane: C...	21.062	3.8	ng/u1	1383760	5	21.386	7350150	20.0
(DEL) Alkane: S...	22.144	2.3	ng/u1	825851	5	21.386	7350150	20.0