

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080924\
 Data File : BM047106.D
 Acq On : 09 Aug 2024 04:51
 Operator : RC/JU
 Sample : P3434-08
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 E05X0

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-BM080224.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BM047106.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.016	18	22	29	rVB	48976	63901	0.48%	0.075%
2	3.275	64	66	81	rVV	79856	132845	0.99%	0.156%
3	3.405	84	88	108	rVB	629668	915453	6.83%	1.078%
4	3.699	135	138	145	rVB	26748	38605	0.29%	0.045%
5	5.528	445	449	453	rBV2	26829	37651	0.28%	0.044%
6	6.593	623	630	641	rVB	945866	1457393	10.88%	1.715%
7	6.745	650	656	670	rVB	709769	1100334	8.21%	1.295%
8	6.928	680	687	700	rVB	925515	1418261	10.59%	1.669%
9	7.028	700	704	712	rBV4	27496	56593	0.42%	0.067%
10	7.387	757	765	772	rBV	1115392	1671774	12.48%	1.968%
11	7.481	775	781	789	rBV5	17986	47943	0.36%	0.056%
12	8.104	881	887	897	rBV	1105630	1713989	12.79%	2.017%
13	8.528	953	959	969	rVV	856223	1401099	10.46%	1.649%
14	9.116	1053	1059	1064	rBV	103134	150791	1.13%	0.177%
15	9.239	1073	1080	1089	rBV	745141	1208325	9.02%	1.422%
16	9.781	1162	1172	1181	rVV	1159221	1915294	14.30%	2.254%
17	10.139	1226	1233	1239	rBV	1426412	2296241	17.14%	2.703%
18	10.192	1239	1242	1247	rVB	35733	52136	0.39%	0.061%
19	10.292	1251	1259	1271	rBV	936412	1631387	12.18%	1.920%
20	10.722	1326	1332	1340	rBV3	69486	186447	1.39%	0.219%
21	10.898	1356	1362	1371	rBV	151445	272309	2.03%	0.321%
22	11.192	1408	1412	1425	rVB5	16061	37327	0.28%	0.044%
23	11.739	1499	1505	1511	rVB3	28399	47980	0.36%	0.056%
24	11.816	1511	1518	1524	rBV	53508	83386	0.62%	0.098%
25	12.039	1552	1556	1564	rVB	44010	68201	0.51%	0.080%
26	12.333	1600	1606	1611	rBV2	23152	36426	0.27%	0.043%
27	12.880	1694	1699	1704	rVB3	27956	43176	0.32%	0.051%
28	12.980	1711	1716	1721	rBV3	18455	33472	0.25%	0.039%
29	13.357	1774	1780	1784	rBV3	36848	61427	0.46%	0.072%
30	13.410	1784	1789	1793	rVV2	28449	44335	0.33%	0.052%
31	13.469	1793	1799	1809	rVV	1895816	2681345	20.02%	3.156%
32	13.551	1809	1813	1817	rVV2	28694	49117	0.37%	0.058%
33	13.598	1817	1821	1824	rVV2	31908	46756	0.35%	0.055%
34	13.727	1836	1843	1855	rVB	2237707	3375729	25.20%	3.973%
35	13.974	1881	1885	1889	rVB2	31335	39067	0.29%	0.046%
36	14.039	1889	1896	1903	rBV	2108593	3035472	22.66%	3.573%

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Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM080224.MA.M
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37	14.157	1910	1916	1924	rVB	112040	157781	1.18%	0.186%
38	14.286	1929	1938	1944	rBV	765625	1321748	9.87%	1.556%
39	14.445	1959	1965	1971	rVB4	27385	56090	0.42%	0.066%
40	14.674	1999	2004	2009	rBV4	20241	39420	0.29%	0.046%
41	14.845	2026	2033	2035	rBV7	20252	37943	0.28%	0.045%
42	14.880	2035	2039	2044	rVB5	26426	39857	0.30%	0.047%
43	14.939	2045	2049	2053	rVB	30900	35568	0.27%	0.042%
44	15.045	2060	2067	2073	rBV	2912157	4032592	30.10%	4.746%
45	15.092	2073	2075	2080	rVB3	41637	56004	0.42%	0.066%
46	15.186	2080	2091	2097	rBV	718221	987457	7.37%	1.162%
47	15.345	2112	2118	2121	rBV5	21836	35452	0.26%	0.042%
48	15.404	2123	2128	2132	rBV	23882	38498	0.29%	0.045%
49	15.498	2139	2144	2148	rBV4	25126	40374	0.30%	0.048%
50	15.545	2149	2152	2160	rVB2	28063	60488	0.45%	0.071%
51	15.633	2161	2167	2174	rVB8	30483	62287	0.46%	0.073%
52	15.804	2191	2196	2197	rBV3	53091	62871	0.47%	0.074%
53	15.827	2197	2200	2207	rVB	93390	136622	1.02%	0.161%
54	15.963	2219	2223	2229	rBV2	49990	74766	0.56%	0.088%
55	16.033	2231	2235	2241	rVV6	17525	33670	0.25%	0.040%
56	16.245	2262	2271	2278	rBV6	47548	99054	0.74%	0.117%
57	16.351	2283	2289	2293	rBV6	27165	60056	0.45%	0.071%
58	16.457	2304	2307	2314	rVB3	26710	48031	0.36%	0.057%
59	16.545	2318	2322	2326	rBV5	23869	41367	0.31%	0.049%
60	16.604	2326	2332	2338	rVV4	52940	112686	0.84%	0.133%
61	16.798	2360	2365	2370	rVV	2500635	3554419	26.53%	4.184%
62	16.839	2370	2372	2376	rVV	179475	222382	1.66%	0.262%
63	16.898	2376	2382	2394	rVB2	3029370	4277830	31.93%	5.035%
64	17.121	2417	2420	2427	rVB5	23040	38103	0.28%	0.045%
65	17.215	2432	2436	2444	rVB4	46339	90479	0.68%	0.106%
66	17.339	2453	2457	2460	rBV	56538	67610	0.50%	0.080%
67	17.380	2460	2464	2468	rVB4	23722	35697	0.27%	0.042%
68	17.621	2499	2505	2508	rBV4	55664	93164	0.70%	0.110%
69	17.674	2510	2514	2518	rBV	193333	261482	1.95%	0.308%
70	17.739	2518	2525	2533	rBV2	1079338	1607450	12.00%	1.892%
71	17.862	2543	2546	2550	rBV3	37341	54889	0.41%	0.065%
72	17.992	2563	2568	2570	rBV4	77613	122714	0.92%	0.144%
73	18.033	2572	2575	2580	rVB2	114585	161207	1.20%	0.190%
74	18.227	2604	2608	2615	rVB	385967	495083	3.70%	0.583%
75	18.609	2670	2673	2679	rVB4	63868	88890	0.66%	0.105%
76	18.680	2681	2685	2688	rBV	85444	113885	0.85%	0.134%

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77	18.751	2694	2697	2704	rBV3	63541	124827	0.93%	0.147%
78	18.845	2709	2713	2714	rBV	56865	67100	0.50%	0.079%
79	18.880	2714	2719	2722	rBV	260150	387547	2.89%	0.456%
80	18.915	2722	2725	2728	rVV3	65493	72850	0.54%	0.086%
81	19.039	2742	2746	2757	rBV	395300	599726	4.48%	0.706%
82	19.215	2769	2776	2783	rBV	3494018	5147143	38.42%	6.058%
83	19.309	2789	2792	2800	rVB4	64181	118531	0.88%	0.140%
84	19.651	2846	2850	2856	rBV2	138022	181385	1.35%	0.213%
85	19.780	2869	2872	2874	rBV2	95973	109154	0.81%	0.128%
86	19.809	2874	2877	2881	rVB	127953	144852	1.08%	0.170%
87	20.092	2923	2925	2929	rVB2	62274	59171	0.44%	0.070%
88	20.133	2929	2932	2943	rBV4	123959	207772	1.55%	0.245%
89	20.303	2957	2961	2965	rBV3	120789	208523	1.56%	0.245%
90	20.486	2989	2992	2994	rBV3	74422	91886	0.69%	0.108%
91	20.774	3037	3041	3053	rVB	1641795	2128826	15.89%	2.506%
92	21.033	3080	3085	3090	rBV	2770034	3771768	28.16%	4.439%
93	21.256	3120	3123	3128	rBV4	151096	214664	1.60%	0.253%
94	21.792	3207	3214	3225	rVB2	821097	1655825	12.36%	1.949%
95	22.639	3355	3358	3365	rVB	244481	375400	2.80%	0.442%
96	23.015	3418	3422	3427	rBV	351939	604162	4.51%	0.711%
97	23.074	3428	3432	3444	rVB4	319470	725758	5.42%	0.854%
98	23.550	3506	3513	3521	rBV	2039748	4431879	33.08%	5.216%
99	23.738	3537	3545	3552	rBV	1651120	3726923	27.82%	4.387%
100	27.144	4112	4124	4140	rVB3	3472588	13395842	100.00%	15.767%

Sum of corrected areas: 84959467

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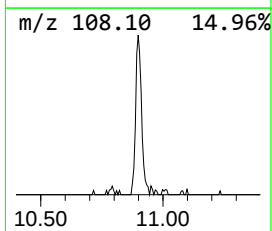
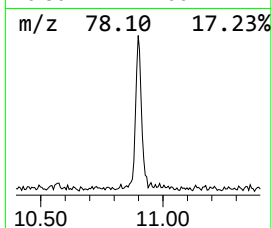
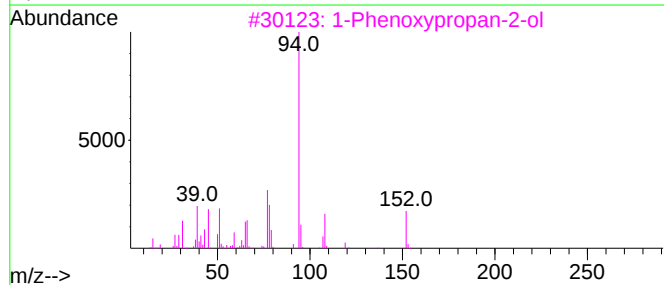
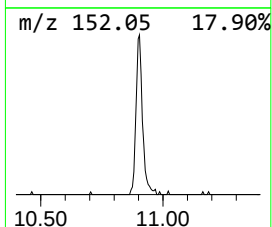
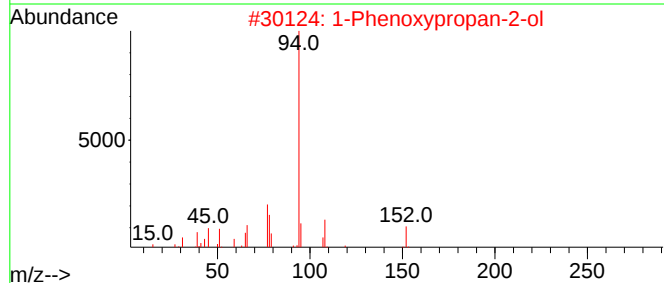
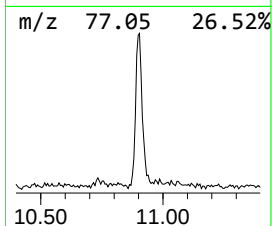
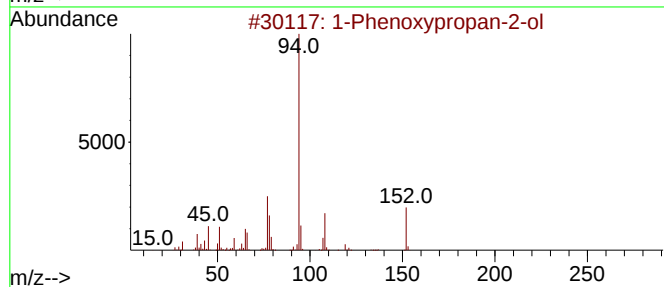
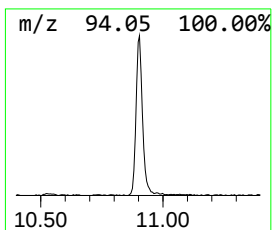
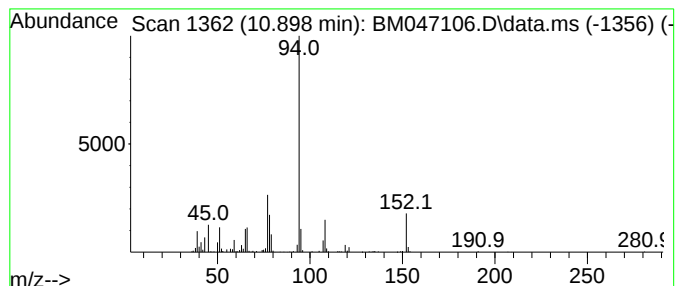
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM080224.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 1-Phenoxypropan-2-ol Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.898	2.37 ng/ul	272309	Naphthalene-d8	10.139

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	96
2			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	95
3			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	94
4			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	87
5			1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	87



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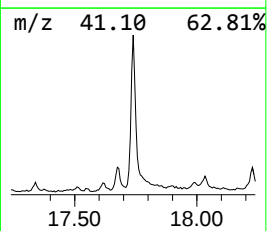
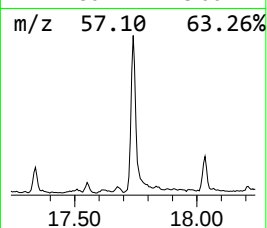
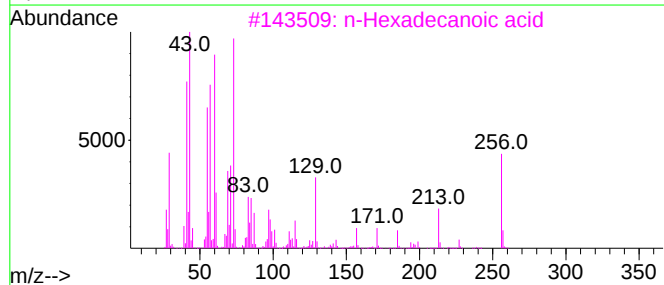
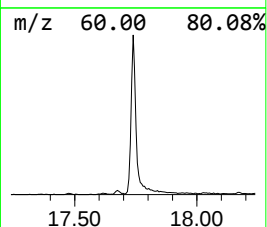
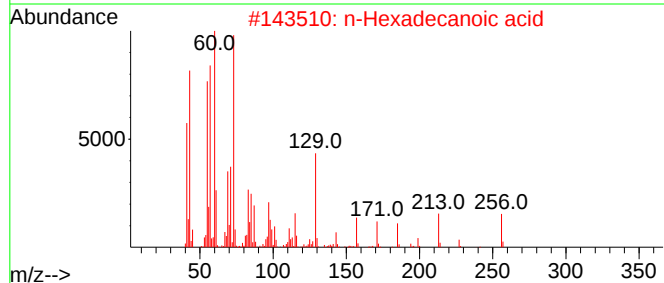
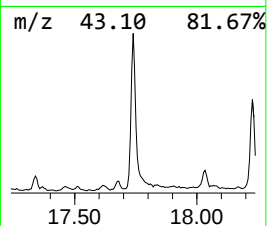
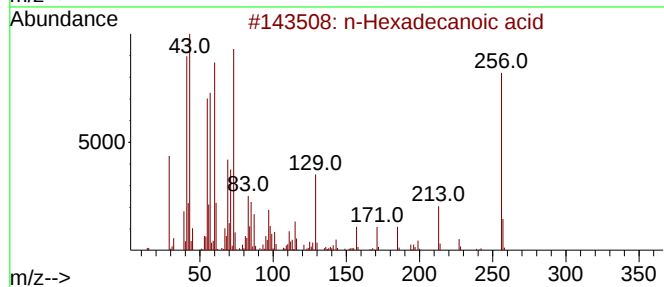
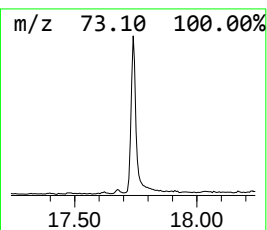
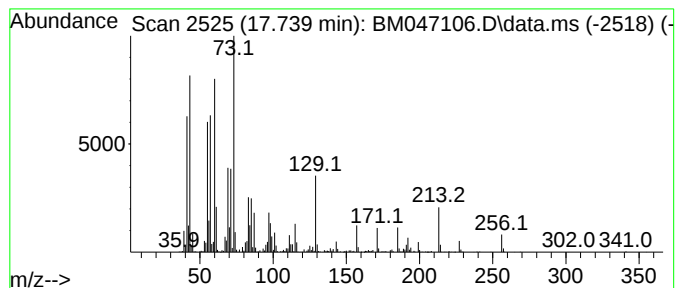
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM080224.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.739	9.04 ng/ul	1607450	Phenanthrene-d10	16.798

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			Tetradecanoic acid	228	C14H28O2	000544-63-8	91



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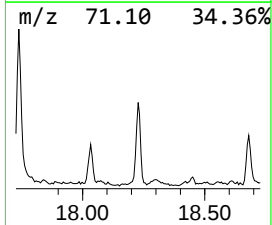
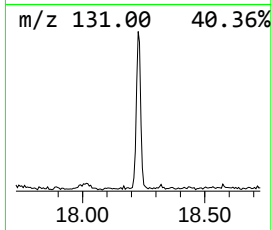
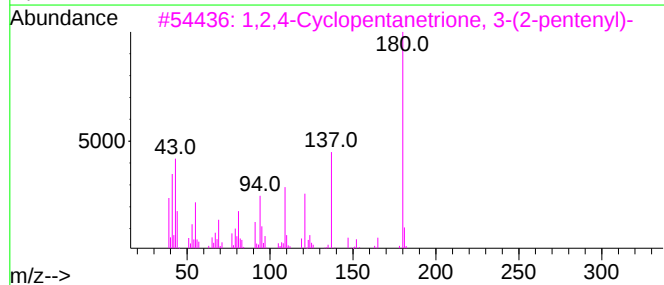
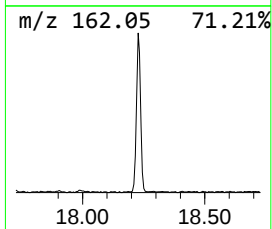
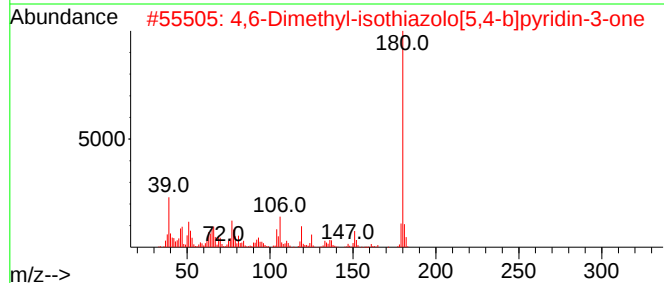
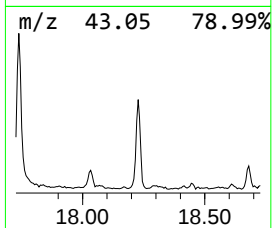
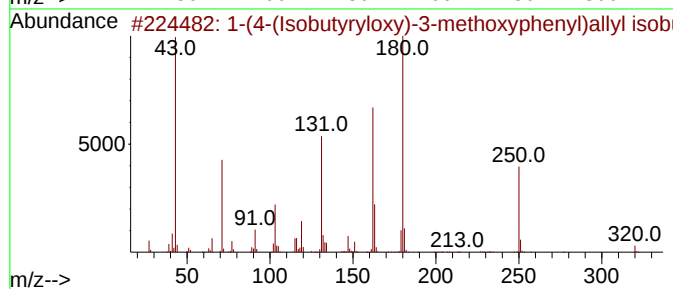
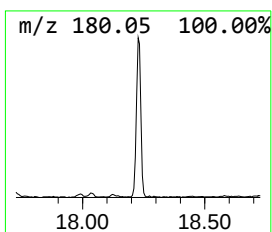
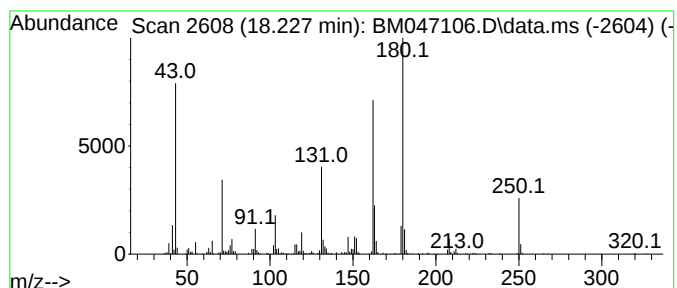
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 1-(4-(Isobutyryloxy)-3-meth... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.227	2.79 ng/ul	495083	Phenanthrene-d10	16.798

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-(4-(Isobutyryloxy)-3-methoxyph...	320	C18H24O5	055249-41-7	76		
2	4,6-Dimethyl-isothiazolo[5,4-b]p...	180	C8H8N2OS	060750-75-6	35		
3	1,2,4-Cyclopentanetrione, 3-(2-p...	180	C10H12O3	054644-27-8	27		
4	Benzaldehyde, m-nitro-, O-methyl...	180	C8H8N2O3	033499-33-1	27		
5	1,2-Dimethyl-3-nitro-4-nitroso-b...	180	C8H8N2O3	1000275-00-6	27		



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Operator : RC/JU
Sample : P3434-08
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
E05X0

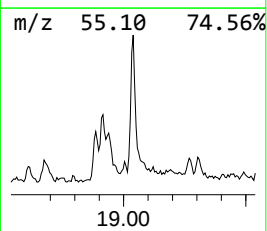
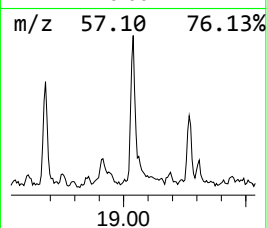
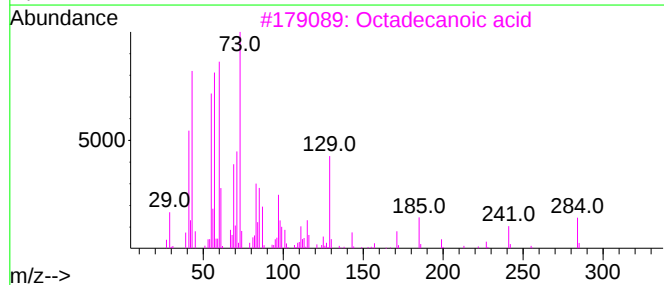
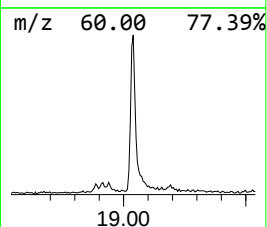
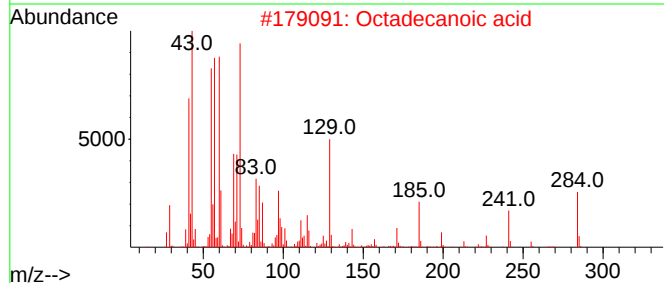
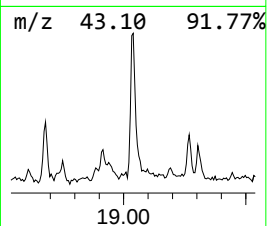
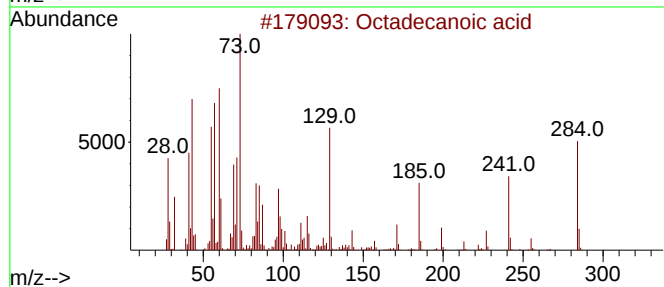
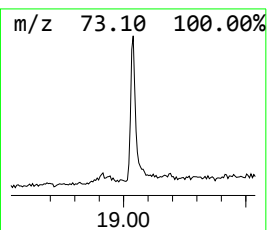
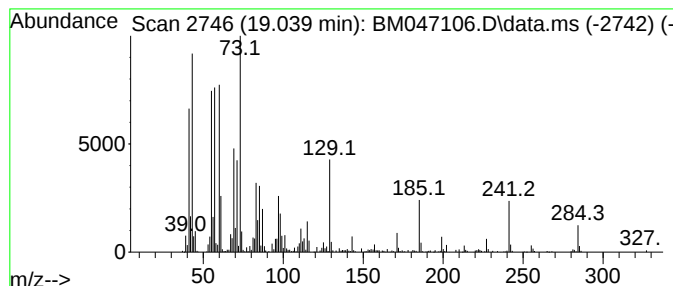
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Octadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.039	3.18 ng/ul	599726	Chrysene-d12	21.033

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid			284	C18H36O2	000057-11-4	99
2	Octadecanoic acid			284	C18H36O2	000057-11-4	98
3	Octadecanoic acid			284	C18H36O2	000057-11-4	98
4	Octadecanoic acid			284	C18H36O2	000057-11-4	95
5	Octadecanoic acid			284	C18H36O2	000057-11-4	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080924\
Data File : BM047106.D
Acq On : 09 Aug 2024 04:51
Operator : RC/JU
Sample : P3434-08
Misc :
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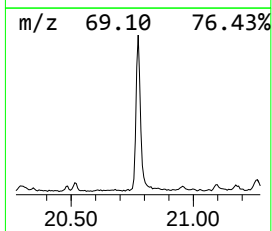
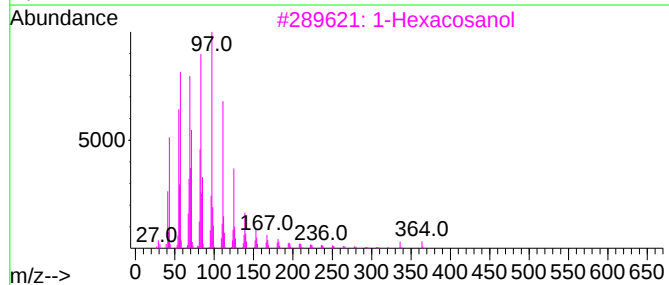
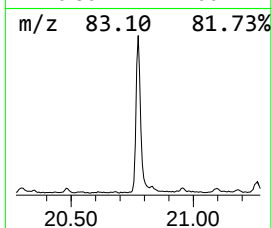
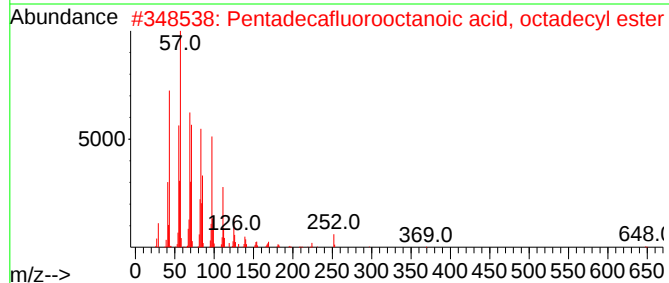
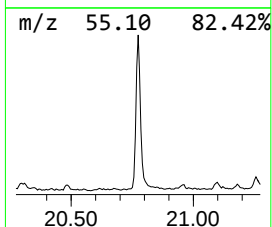
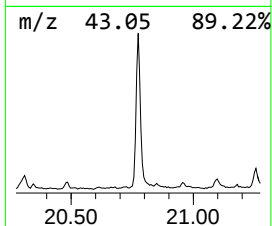
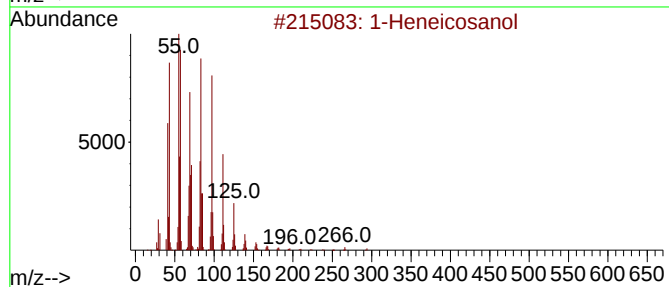
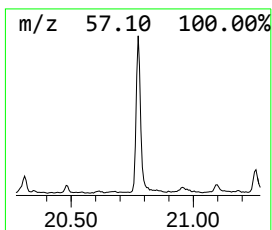
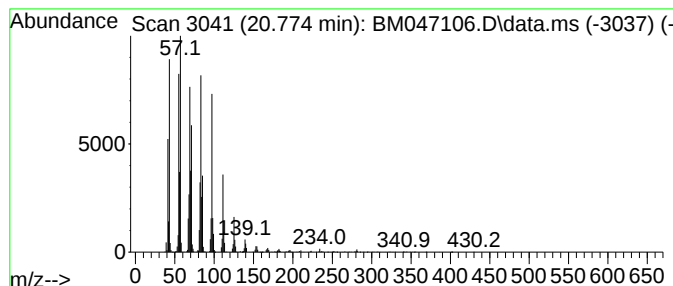
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 1-Heneicosanol Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
20.774	11.29 ng/ul	2128830	Chrysene-d12	21.033	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heneicosanol	312	C21H44O	015594-90-8	94
2	Pentadecafluorooctanoic acid, octadecyl ...	666	C26H37F15O2	1000406-04-8	93
3	1-Hexacosanol	382	C26H54O	000506-52-5	91
4	Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	91
5	Behenic alcohol	326	C22H46O	000661-19-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080924\
Data File : BM047106.D
Acq On : 09 Aug 2024 04:51
Operator : RC/JU
Sample : P3434-08
Misc :
ALS Vial : 23 Sample Multiplier: 1

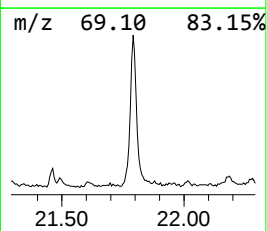
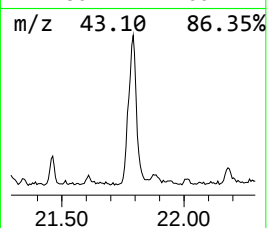
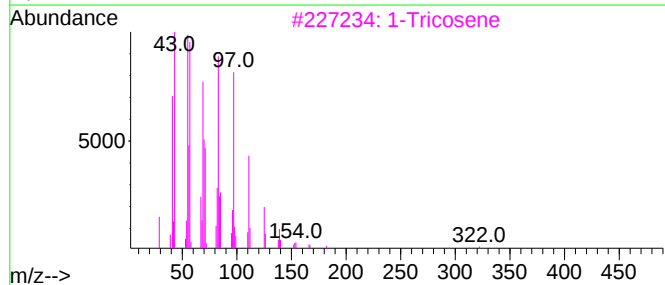
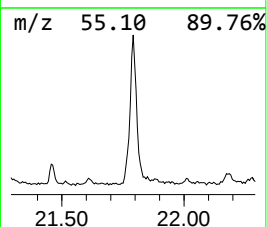
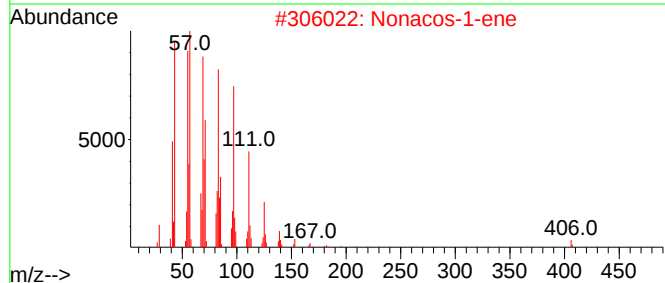
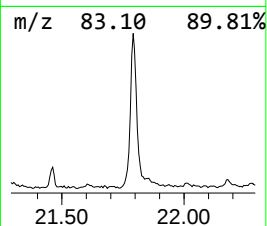
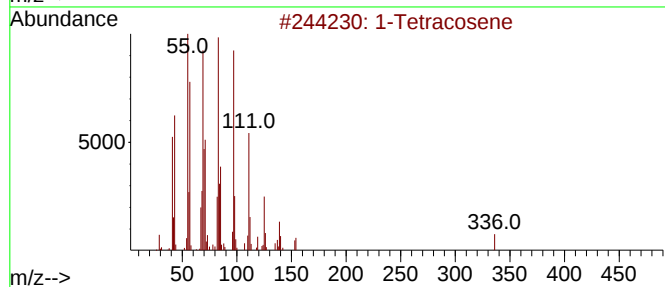
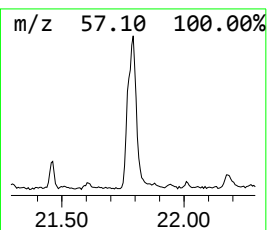
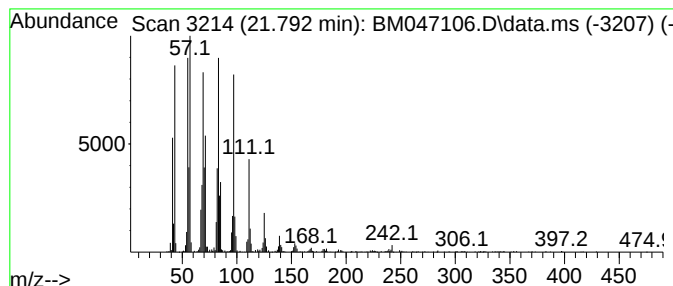
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD		R.T.	
21.792	8.78 ng/ul	1655830	Chrysene-d12		21.033	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	1-Tetracosene		336	C24H48	010192-32-2	99
2	Nonacos-1-ene		406	C29H58	018835-35-3	95
3	1-Tricosene		322	C23H46	018835-32-0	95
4	1-Heptacosanol		396	C27H56O	002004-39-9	94
5	Behenic alcohol		326	C22H46O	000661-19-8	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080924\
Data File : BM047106.D
Acq On : 09 Aug 2024 04:51
Operator : RC/JU
Sample : P3434-08
Misc :
ALS Vial : 23 Sample Multiplier: 1

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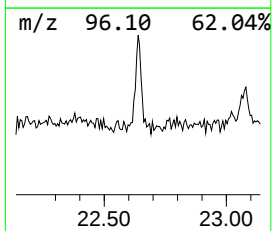
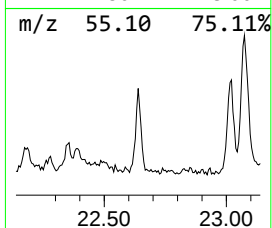
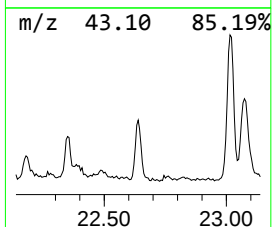
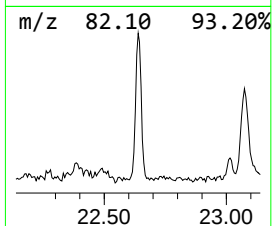
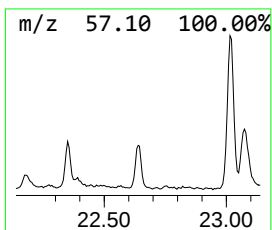
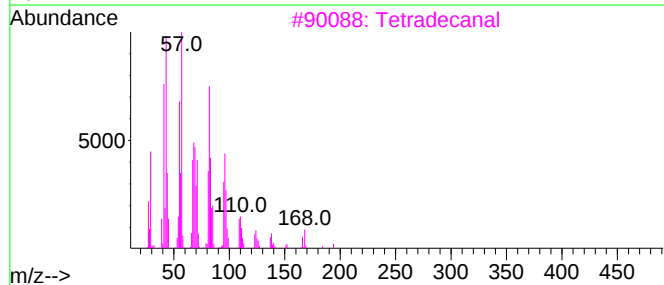
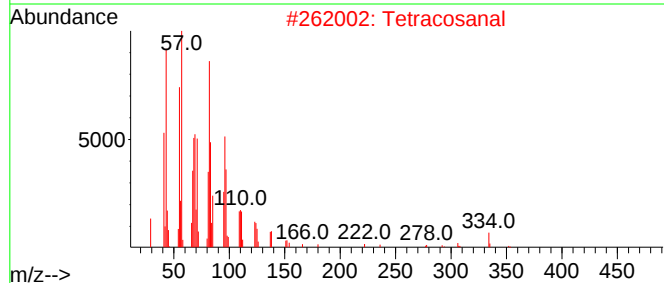
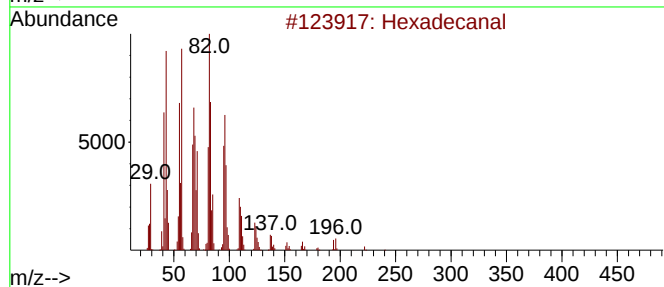
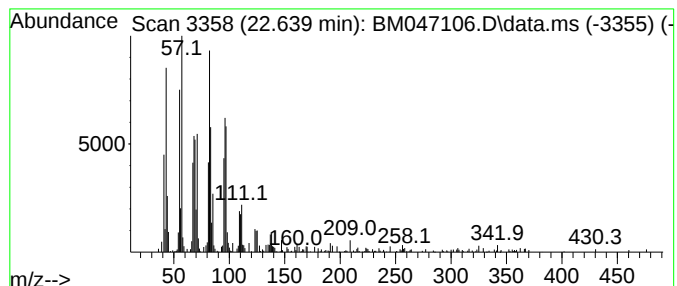
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Hexadecanal Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.639	2.01 ng/ul	375400	Perylene-d12	23.738

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecanal	240	C16H32O	000629-80-1	98
2			Tetracosanal	352	C24H48O	057866-08-7	95
3			Tetradecanal	212	C14H28O	000124-25-4	94
4			Docosanal	324	C22H44O	057402-36-5	94
5			Tetradecanal	212	C14H28O	000124-25-4	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080924\
Data File : BM047106.D
Acq On : 09 Aug 2024 04:51
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Sample : P3434-08
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ClientSampleId :
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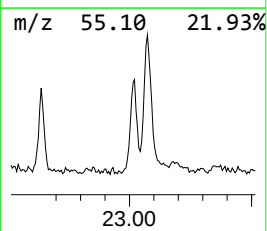
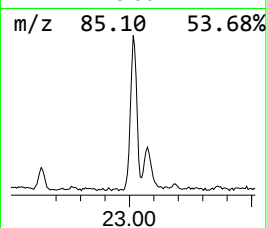
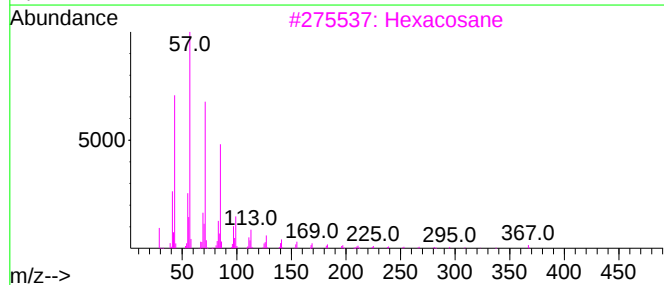
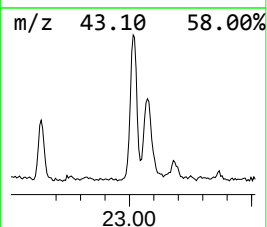
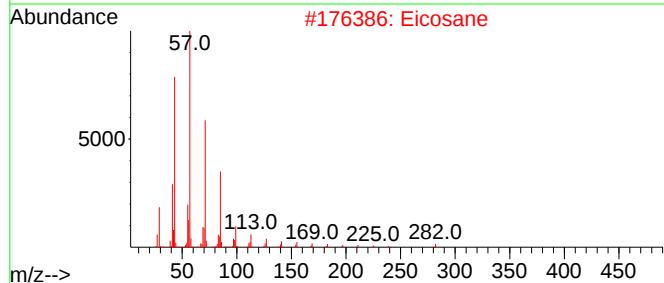
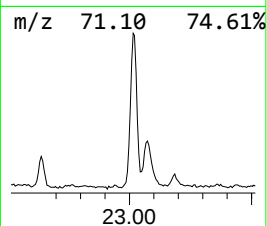
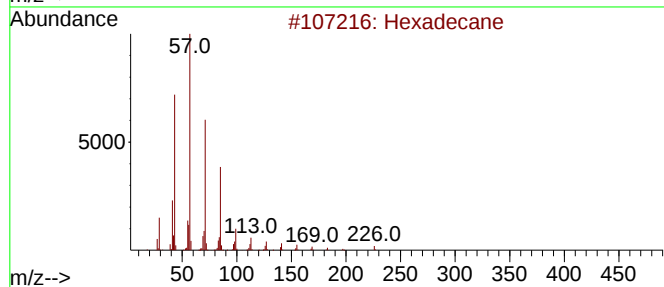
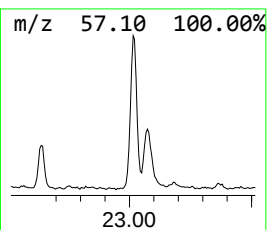
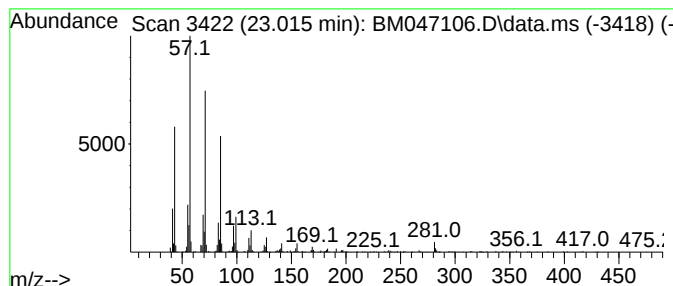
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.015	3.24 ng/ul	604162	Perylene-d12	23.738

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecane	226	C16H34	000544-76-3	93
2			Eicosane	282	C20H42	000112-95-8	93
3			Hexacosane	366	C26H54	000630-01-3	90
4			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	90
5			Heneicosane	296	C21H44	000629-94-7	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080924\
Data File : BM047106.D
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Sample : P3434-08
Misc :
ALS Vial : 23 Sample Multiplier: 1

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ClientSampleId :
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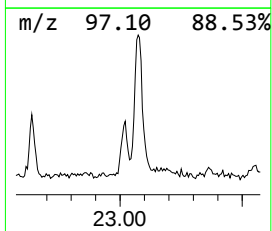
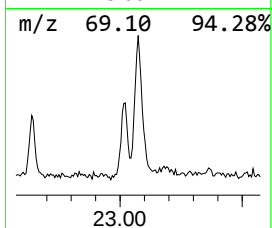
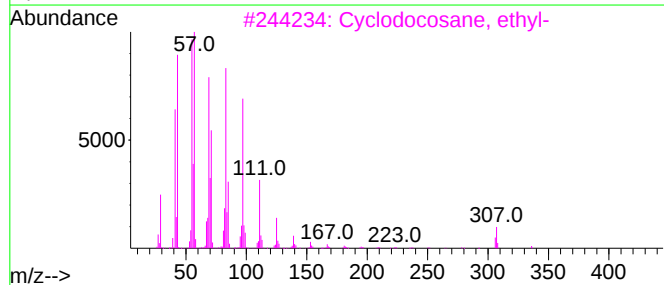
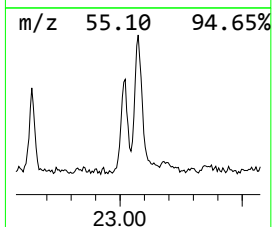
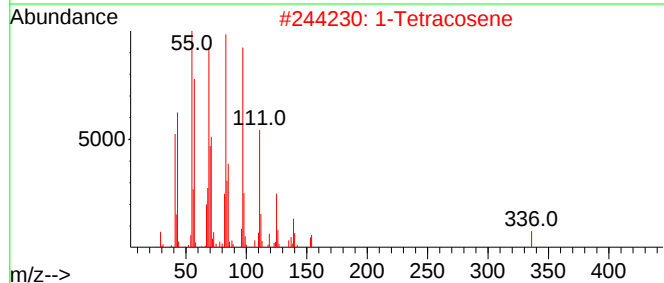
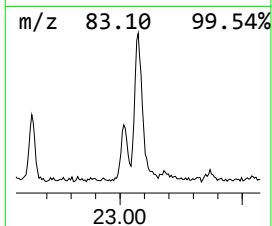
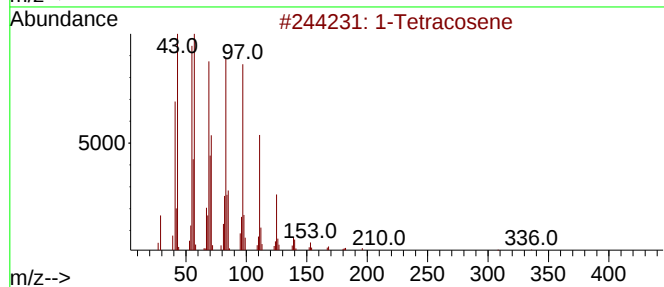
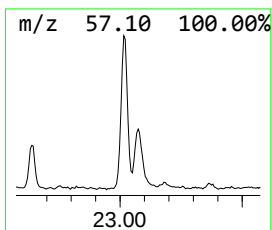
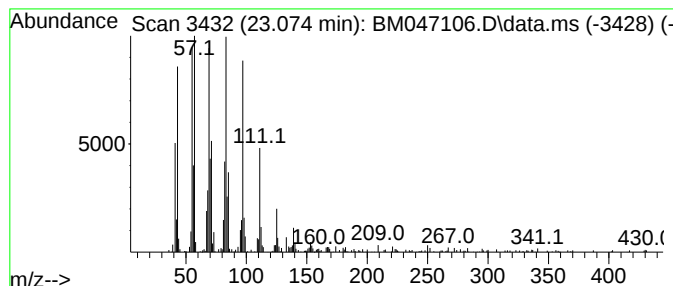
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.074	3.89 ng/ul	725758	Perylene-d12	23.738

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Tetracosene	336	C24H48	010192-32-2	99
2		1-Tetracosene	336	C24H48	010192-32-2	98
3		Cyclodocosane, ethyl-	336	C24H48	1000151-22-6	94
4		Pentacos-1-ene	350	C25H50	016980-85-1	94
5		1-Hexacosene	364	C26H52	018835-33-1	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080924\
Data File : BM047106.D
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Sample : P3434-08
Misc :
ALS Vial : 23 Sample Multiplier: 1

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ClientSampleId :
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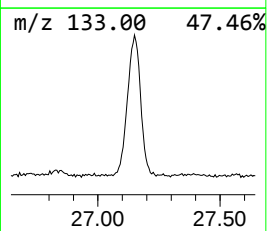
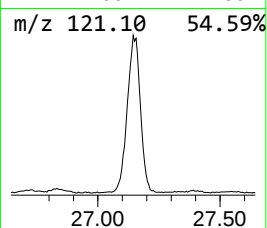
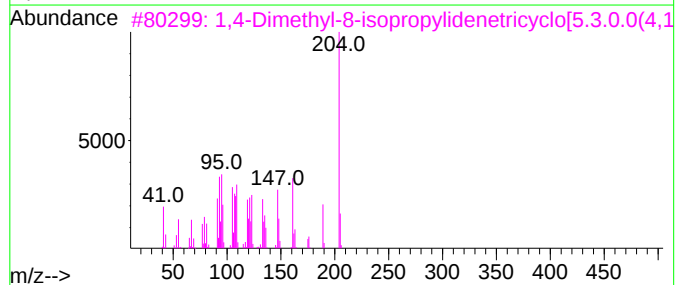
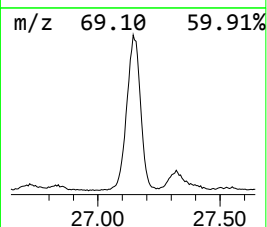
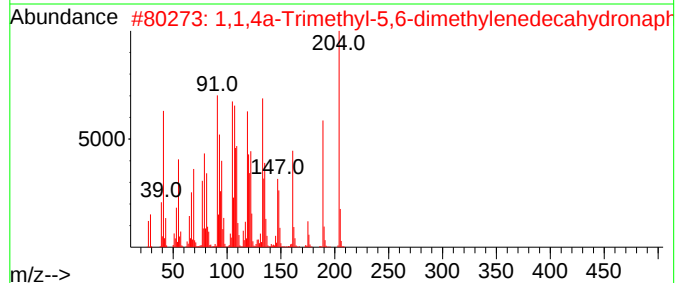
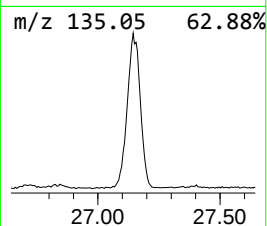
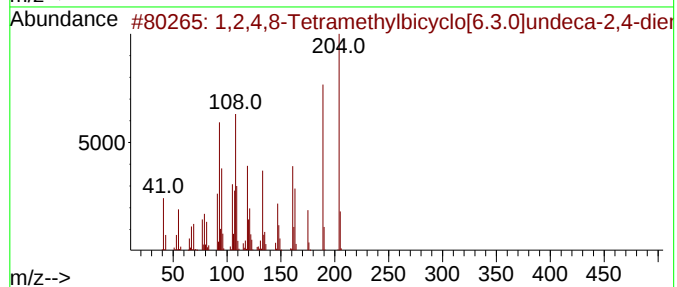
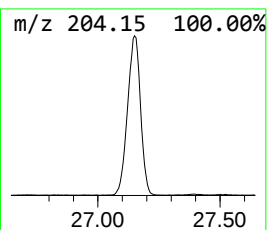
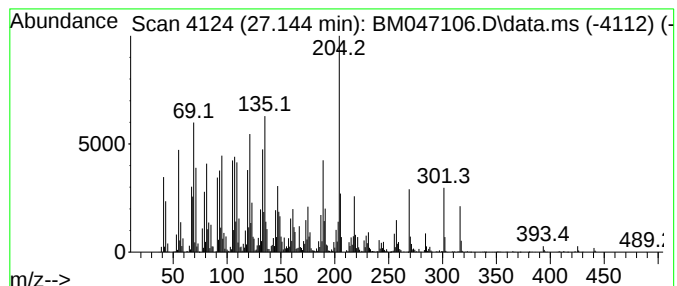
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 1,2,4,8-Tetramethylbicyclo[... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
27.144	71.89 ng/ul	13395800	Perylene-d12	23.738

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,2,4,8-Tetramethylbicyclo[6.3.0...	204	C15H24	137235-51-9	70		
2	1,1,4a-Trimethyl-5,6-dimethylene...	204	C15H24	1000193-60-8	53		
3	1,4-Dimethyl-8-isopropylidenetri...	204	C15H24	1010140-07-7	50		
4	1H-3a,7-Methanoazulene, 2,3,6,7,...	204	C15H24	000560-32-7	49		
5	2H-Cyclopentacyclooctene, 4,5,6,...	204	C15H24	1000221-85-8	43		



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080924\
Data File : BM047106.D
Acq On : 09 Aug 2024 04:51
Operator : RC/JU
Sample : P3434-08
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
E05X0

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM080224.MA.M
Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
1-Phenoxypropan...	10.898	2.4	ng/ul	272309	2	10.139	2296240	20.0
n-Hexadecanoic ...	17.739	9.0	ng/ul	1607450	4	16.798	3554420	20.0
1-(4-(Isobutyry...	18.227	2.8	ng/ul	495083	4	16.798	3554420	20.0
Octadecanoic acid	19.039	3.2	ng/ul	599726	5	21.033	3771770	20.0
1-Heneicosanol	20.774	11.3	ng/ul	2128830	5	21.033	3771770	20.0
(DEL) Alkane: S...	21.792	8.8	ng/ul	1655830	5	21.033	3771770	20.0
Hexadecanal	22.639	2.0	ng/ul	375400	6	23.738	3726920	20.0
(DEL) Alkane: S...	23.015	3.2	ng/ul	604162	6	23.738	3726920	20.0
(DEL) Alkane: S...	23.074	3.9	ng/ul	725758	6	23.738	3726920	20.0
1,2,4,8-Tetrame...	27.144	71.9	ng/ul	13395800	6	23.738	3726920	20.0