

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040822\
 Data File : BP009727.D
 Acq On : 08 Apr 2022 21:01
 Operator : CG/JU
 Sample : N2182-08
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0R15

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_P\Methods\SFAM-EPA-BP040722.M
 Title : SVOA CALIBRATION

Signal : TIC: BP009727.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.128	3	7	24	rVB	604262	889652	36.60%	3.356%
2	3.393	48	52	58	rVB4	9020	13982	0.58%	0.053%
3	3.817	120	124	134	rBV	33629	58507	2.41%	0.221%
4	3.917	138	141	147	rVB2	9193	12708	0.52%	0.048%
5	4.146	175	180	184	rBV3	7545	11741	0.48%	0.044%
6	4.364	212	217	222	rBV4	8467	13509	0.56%	0.051%
7	4.781	283	288	294	rVB6	5871	8871	0.36%	0.033%
8	5.070	333	337	345	rVV4	9619	16843	0.69%	0.064%
9	7.122	679	686	697	rVB	191878	330534	13.60%	1.247%
10	7.299	707	716	727	rBV	153741	253354	10.42%	0.956%
11	7.499	744	750	760	rBV	190506	314400	12.94%	1.186%
12	7.611	765	769	775	rVB5	4175	7903	0.33%	0.030%
13	7.975	823	831	841	rBV	383279	635477	26.15%	2.397%
14	8.675	942	950	959	rBV	206943	340404	14.01%	1.284%
15	8.799	964	971	979	rVV2	37516	66986	2.76%	0.253%
16	9.128	1013	1027	1037	rBV	201138	347388	14.29%	1.311%
17	9.605	1101	1108	1113	rVB4	6539	11193	0.46%	0.042%
18	9.857	1143	1151	1163	rBV	154589	264594	10.89%	0.998%
19	9.987	1168	1173	1180	rVV4	7009	11878	0.49%	0.045%
20	10.399	1235	1243	1253	rVB	265203	452314	18.61%	1.706%
21	10.781	1300	1308	1316	rBV	550867	968453	39.85%	3.653%
22	10.910	1323	1330	1341	rBV	147403	261492	10.76%	0.986%
23	11.316	1392	1399	1404	rBV10	4074	9671	0.40%	0.036%
24	11.569	1436	1442	1448	rBV6	5690	11115	0.46%	0.042%
25	12.222	1547	1553	1559	rBV6	5913	12501	0.51%	0.047%
26	12.363	1570	1577	1583	rVV6	6985	13305	0.55%	0.050%
27	12.440	1583	1590	1595	rVV3	12567	20672	0.85%	0.078%
28	12.657	1620	1627	1634	rVB	10257	16959	0.70%	0.064%
29	12.904	1664	1669	1676	rBV	50365	84050	3.46%	0.317%
30	13.040	1687	1692	1696	rBV8	5512	9298	0.38%	0.035%
31	13.451	1756	1762	1767	rBV2	19060	27880	1.15%	0.105%
32	13.928	1838	1843	1848	rVV2	9199	15783	0.65%	0.060%
33	14.010	1848	1857	1864	rVV	628662	869919	35.79%	3.282%
34	14.057	1864	1865	1871	rVV6	9685	14148	0.58%	0.053%
35	14.110	1871	1874	1878	rVV5	6537	9724	0.40%	0.037%
36	14.157	1878	1882	1887	rVV7	5341	12227	0.50%	0.046%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
 Title : SVOA CALIBRATION

37	14.298	1897	1906	1917	rBV	577118	872635	35.90%	3.292%
38	14.522	1939	1944	1949	rBV2	8347	13168	0.54%	0.050%
39	14.604	1950	1958	1965	rVV	917101	1396074	57.44%	5.267%
40	14.669	1965	1969	1977	rVB	21284	36274	1.49%	0.137%
41	14.775	1981	1987	1998	rBV	265864	395363	16.27%	1.492%
42	14.934	2009	2014	2020	rBV4	12847	19571	0.81%	0.074%
43	15.016	2021	2028	2034	rBV4	34175	65423	2.69%	0.247%
44	15.234	2059	2065	2070	rVV10	5101	12669	0.52%	0.048%
45	15.469	2102	2105	2111	rVB3	7051	8961	0.37%	0.034%
46	15.598	2120	2127	2133	rBV	849438	1213894	49.95%	4.579%
47	15.651	2133	2136	2140	rVV3	26353	34997	1.44%	0.132%
48	15.704	2140	2145	2152	rVV	240485	331151	13.63%	1.249%
49	15.775	2153	2157	2161	rVB4	9616	13687	0.56%	0.052%
50	15.840	2164	2168	2171	rVV5	5962	8112	0.33%	0.031%
51	15.951	2183	2187	2193	rVB5	9785	18642	0.77%	0.070%
52	16.098	2207	2212	2218	rVB2	9611	17957	0.74%	0.068%
53	16.334	2246	2252	2254	rBV7	8460	14185	0.58%	0.054%
54	16.363	2254	2257	2261	rVB5	7676	9879	0.41%	0.037%
55	16.475	2271	2276	2278	rBV5	5408	9596	0.39%	0.036%
56	16.757	2321	2324	2330	rVB7	12980	23049	0.95%	0.087%
57	16.892	2344	2347	2352	rBV7	8011	11192	0.46%	0.042%
58	17.004	2362	2366	2371	rVB	16018	22121	0.91%	0.083%
59	17.063	2372	2376	2378	rBV4	7515	8385	0.35%	0.032%
60	17.139	2385	2389	2392	rVB4	9208	11683	0.48%	0.044%
61	17.175	2392	2395	2401	rVB	15898	24111	0.99%	0.091%
62	17.257	2405	2409	2416	rBV2	37035	50010	2.06%	0.189%
63	17.357	2416	2426	2430	rBV	1213916	1797718	73.97%	6.782%
64	17.398	2430	2433	2437	rVB	234716	307862	12.67%	1.161%
65	17.457	2437	2443	2448	rBV	966027	1392439	57.29%	5.253%
66	17.545	2454	2458	2464	rVB7	15032	24405	1.00%	0.092%
67	17.751	2486	2493	2499	rBV2	35155	64511	2.65%	0.243%
68	17.875	2509	2514	2517	rBV5	10995	14977	0.62%	0.057%
69	18.034	2538	2541	2544	rVB5	7654	8939	0.37%	0.034%
70	18.128	2551	2557	2563	rBV5	40495	77318	3.18%	0.292%
71	18.186	2563	2567	2570	rVV	39583	51766	2.13%	0.195%
72	18.245	2570	2577	2585	rVB2	112003	221999	9.13%	0.837%
73	18.345	2591	2594	2599	rVB4	10222	16342	0.67%	0.062%
74	18.410	2600	2605	2608	rBV2	40587	63988	2.63%	0.241%
75	18.534	2622	2626	2632	rBV5	40158	67248	2.77%	0.254%
76	18.704	2651	2655	2657	rBV	22557	28318	1.17%	0.107%

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 Title : SVOA CALIBRATION

77	18.733	2657	2660	2665	rVB2	18274	26271	1.08%	0.099%
78	18.786	2667	2669	2673	rVV3	7372	9477	0.39%	0.036%
79	19.098	2717	2722	2727	rVB4	36437	54517	2.24%	0.206%
80	19.145	2727	2730	2733	rBV5	14958	21411	0.88%	0.081%
81	19.233	2742	2745	2747	rBV2	23340	29107	1.20%	0.110%
82	19.310	2755	2758	2760	rBV2	21799	30244	1.24%	0.114%
83	19.357	2761	2766	2775	rVB	432880	572554	23.56%	2.160%
84	19.475	2783	2786	2793	rVB6	31356	43324	1.78%	0.163%
85	19.681	2815	2821	2825	rBV	1494873	2171459	89.35%	8.192%
86	19.881	2853	2855	2859	rVB2	19440	22805	0.94%	0.086%
87	19.945	2862	2866	2869	rBV2	16539	25210	1.04%	0.095%
88	20.033	2875	2881	2886	rBV4	24004	61945	2.55%	0.234%
89	20.192	2904	2908	2911	rVB2	41753	56125	2.31%	0.212%
90	20.292	2921	2925	2928	rBV3	28582	34811	1.43%	0.131%
91	21.151	3067	3071	3077	rVB3	223792	307650	12.66%	1.161%
92	21.439	3113	3120	3124	rBV	1562226	2430414	100.00%	9.169%
93	21.475	3124	3126	3133	rVB	155826	182267	7.50%	0.688%
94	22.092	3227	3231	3239	rVB2	165244	264327	10.88%	0.997%
95	23.151	3408	3411	3412	rBV2	89039	91193	3.75%	0.344%
96	23.227	3421	3424	3432	rVB2	200233	323813	13.32%	1.222%
97	23.751	3507	3513	3519	rBV2	922132	1731261	71.23%	6.531%
98	23.910	3534	3540	3558	rVB2	1006714	2187686	90.01%	8.253%
99	24.616	3655	3660	3667	rVB	152560	316721	13.03%	1.195%
100	24.698	3670	3674	3682	rVB4	152473	313069	12.88%	1.181%

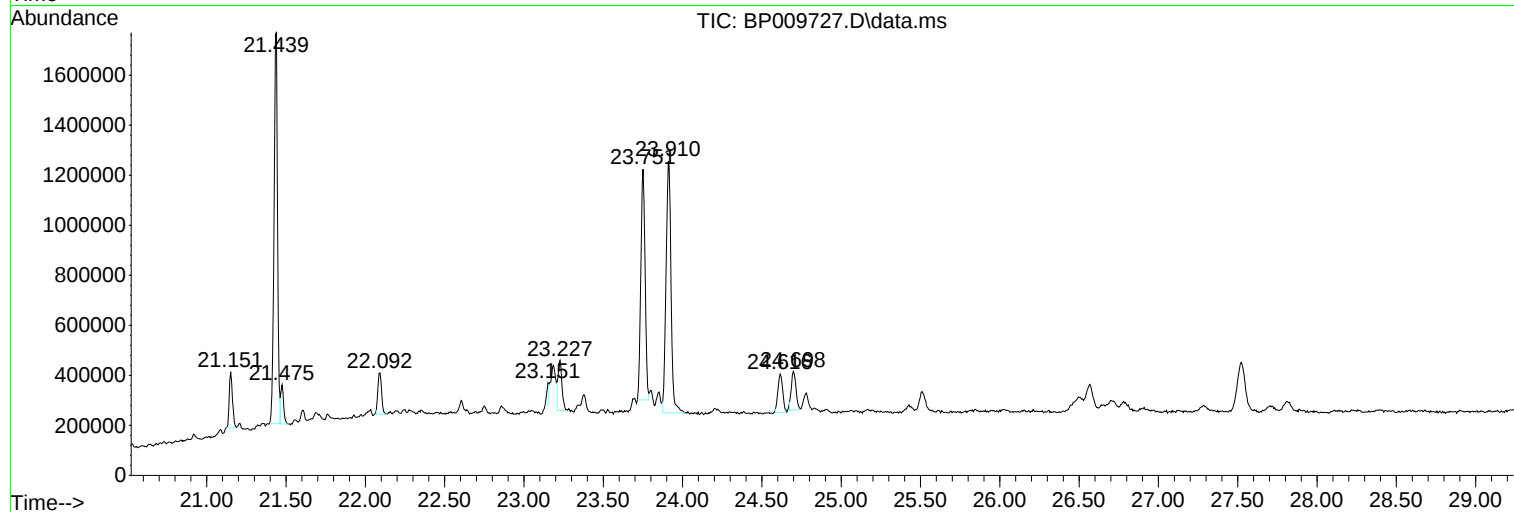
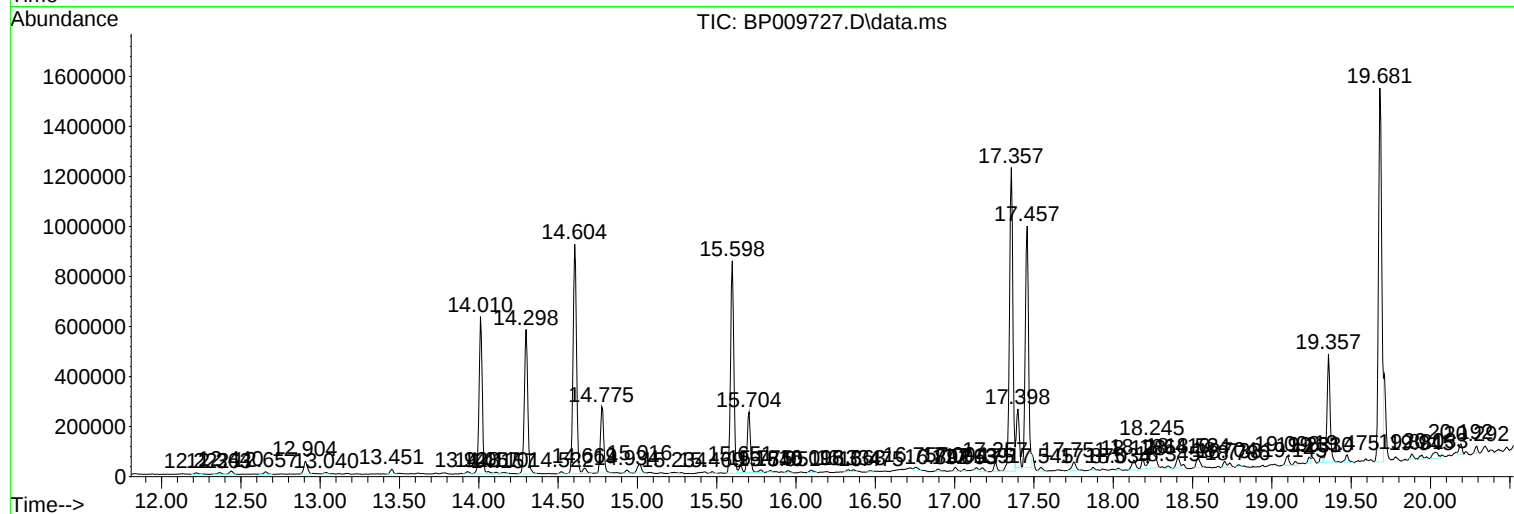
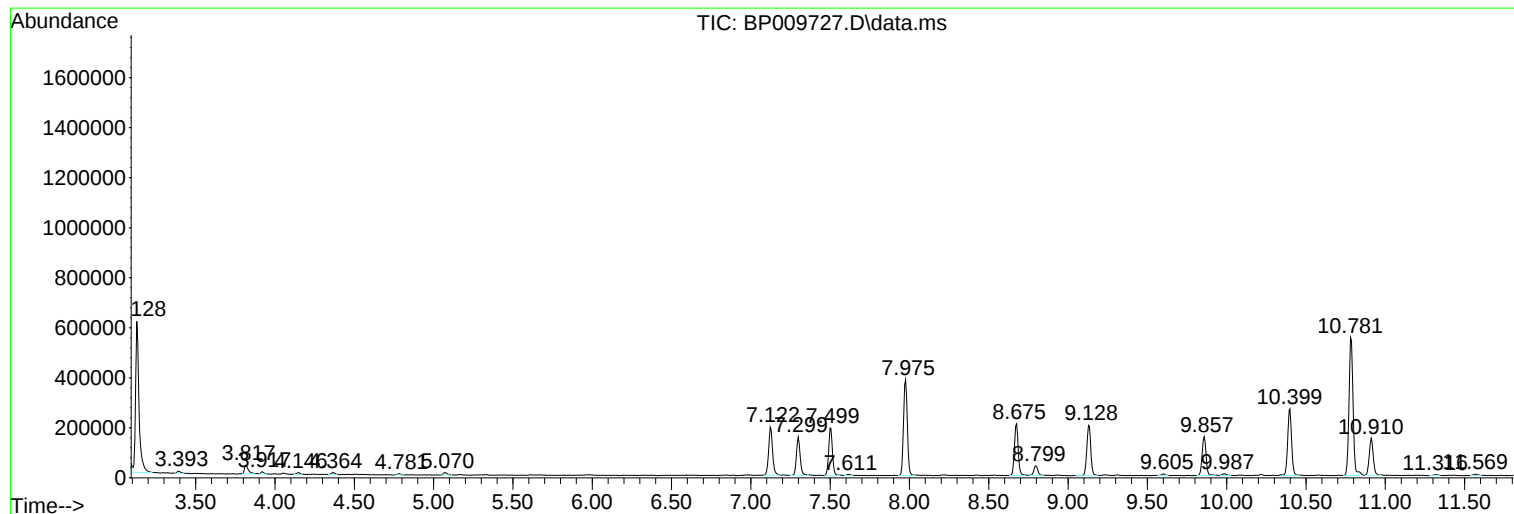
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION

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



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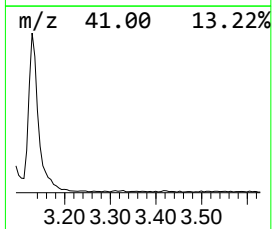
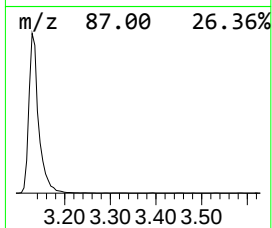
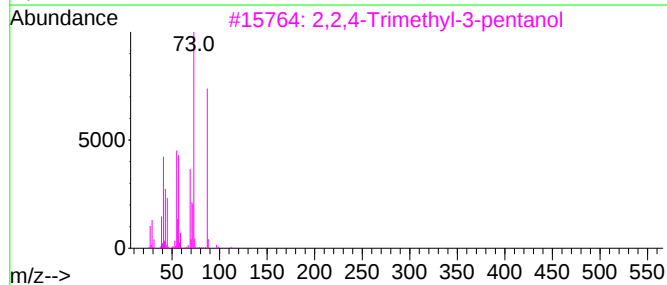
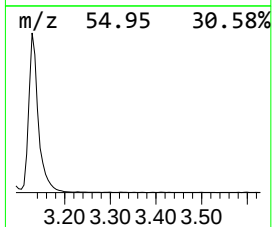
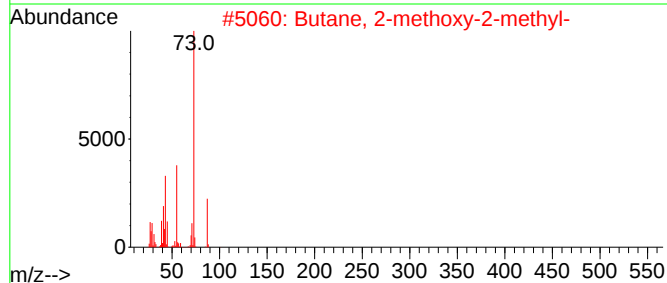
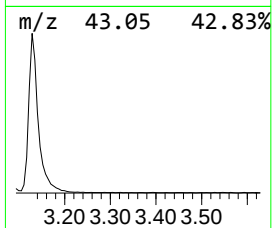
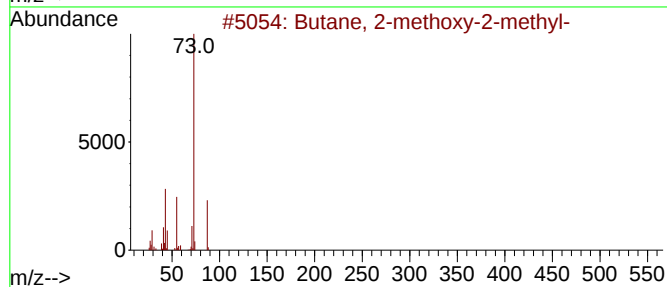
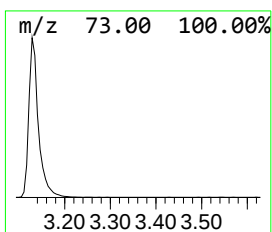
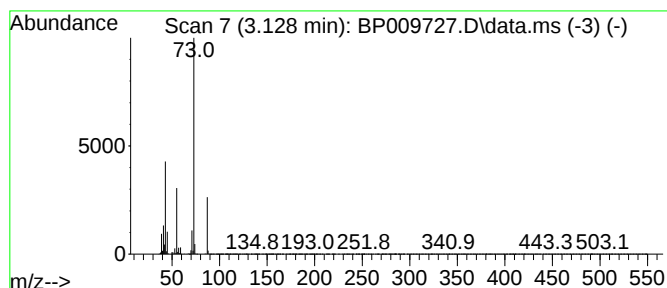
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.128	28.00 ng/ul	889652	1,4-Dichlorobenzene-d4	7.975

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83		
2	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78		
3	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40		
4	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40		
5	1,3-Dioxolane, 2-propyl-	116	C6H12O2	003390-13-4	17		



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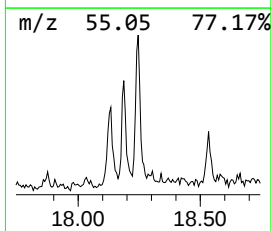
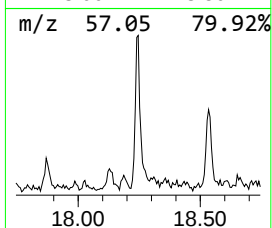
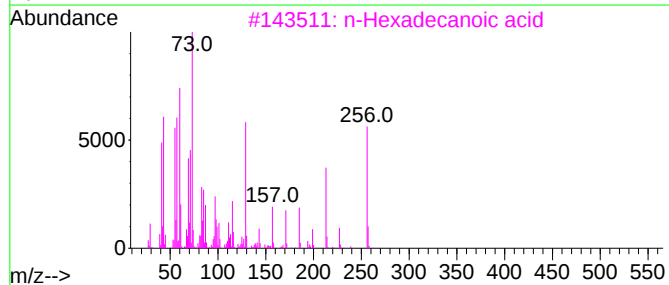
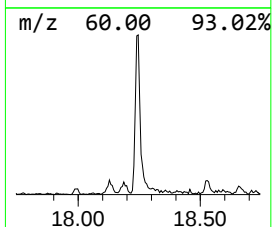
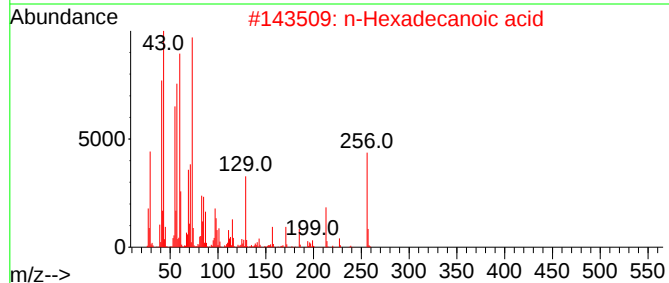
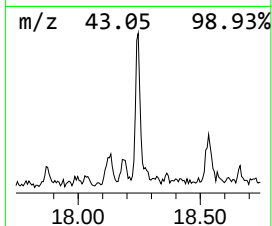
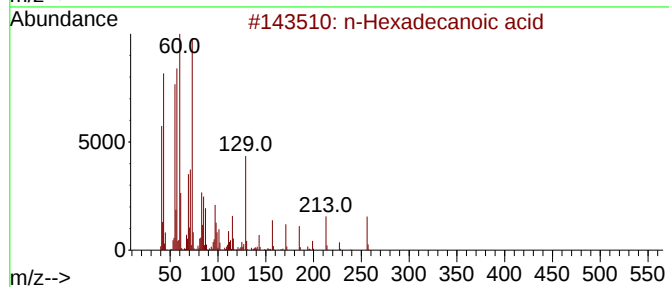
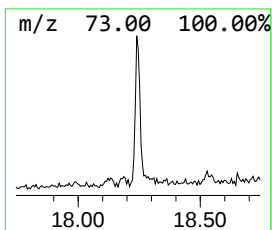
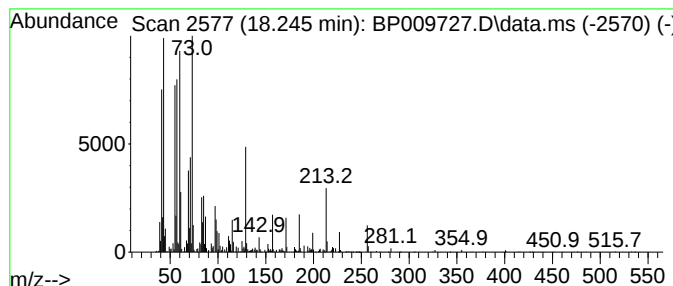
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 n-Hexadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.245	2.47 ng/ul	221999	Phenanthrene-d10	17.357

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	97
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
5			Pentadecanoic acid	242	C15H30O2	001002-84-2	87



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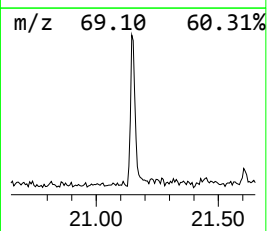
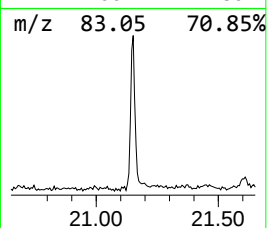
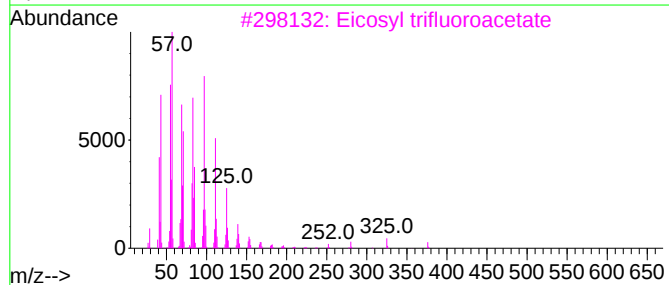
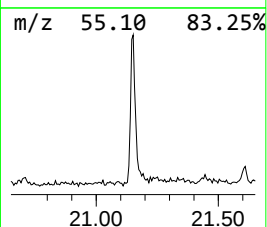
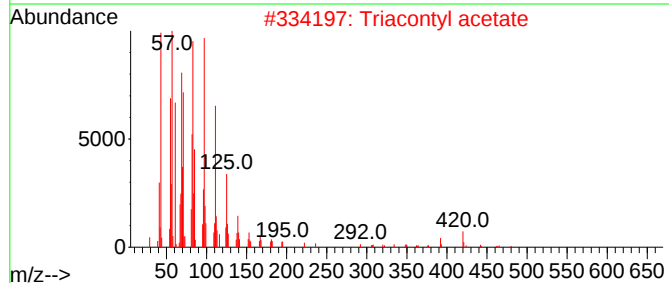
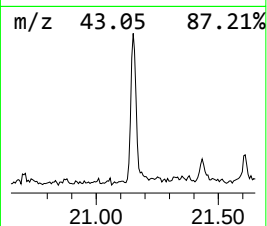
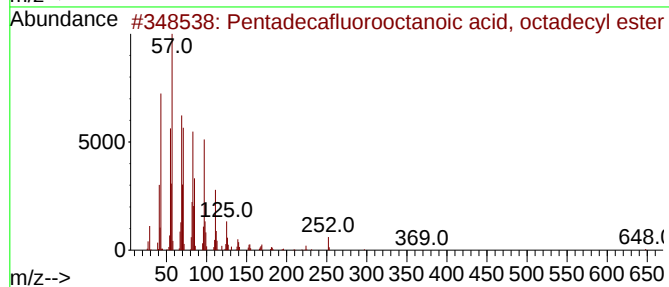
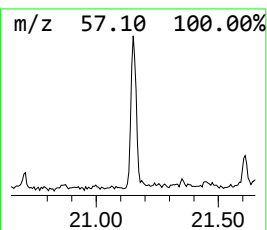
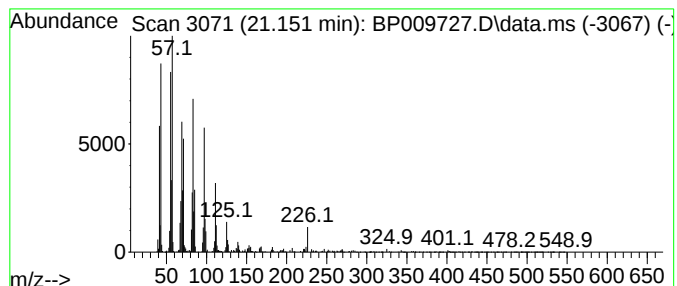
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Pentadecafluorooctanoic aci... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.151	2.53 ng/ul	307650	Chrysene-d12	21.439

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	94
2			Triacontyl acetate	480	C32H64O2	041755-58-2	93
3			Eicosyl trifluoroacetate	394	C22H41F3O2	1000351-75-2	91
4			Pentafluoropropionic acid, octad...	416	C21H37F5O2	959261-25-7	91
5			Nonadecyl pentafluoropropionate	430	C22H39F5O2	1000351-88-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040822\
Data File : BP009727.D
Acq On : 08 Apr 2022 21:01
Operator : CG/JU
Sample : N2182-08
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0R15

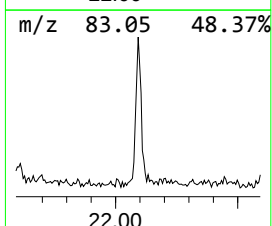
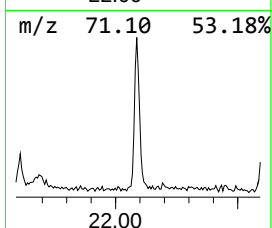
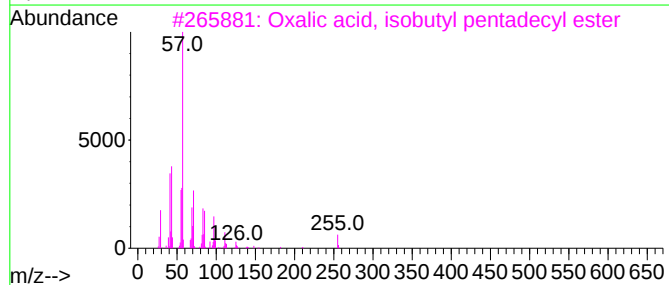
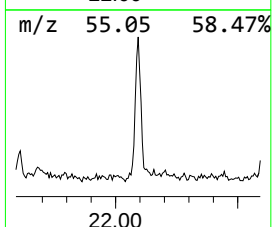
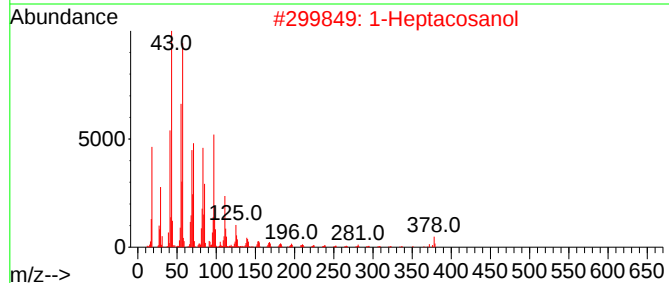
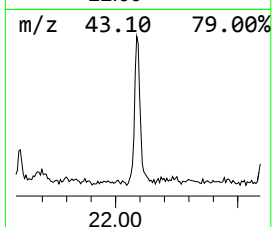
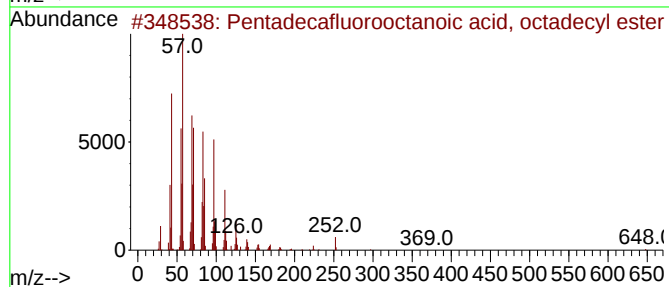
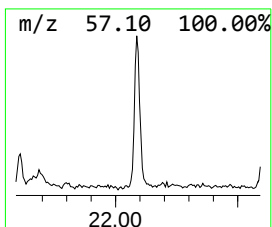
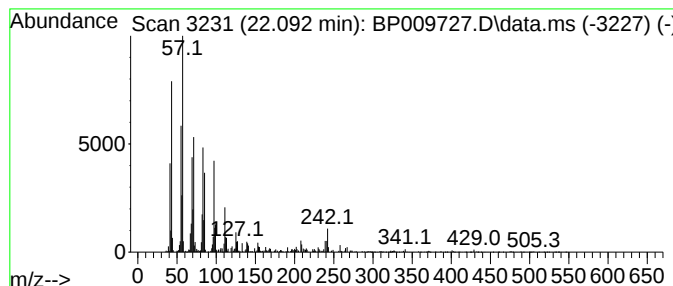
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 1-Heptacosanol Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.092	2.18 ng/ul	264327	Chrysene-d12	21.439

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	87
2			1-Heptacosanol	396	C27H56O	002004-39-9	87
3			Oxalic acid, isobutyl pentadecyl...	356	C21H40O4	1000309-38-0	87
4			9-Tricosene, (Z)-	322	C23H46	027519-02-4	86
5			Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040822\
Data File : BP009727.D
Acq On : 08 Apr 2022 21:01
Operator : CG/JU
Sample : N2182-08
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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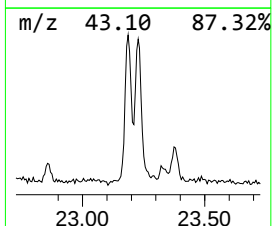
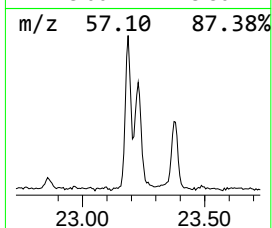
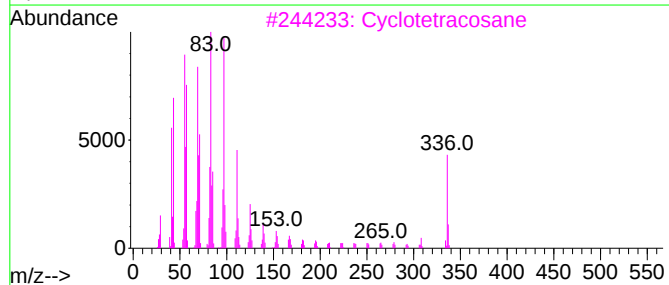
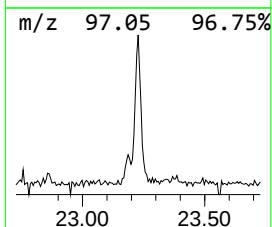
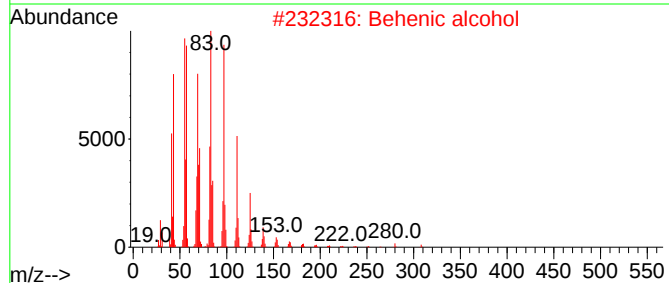
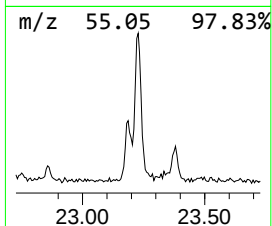
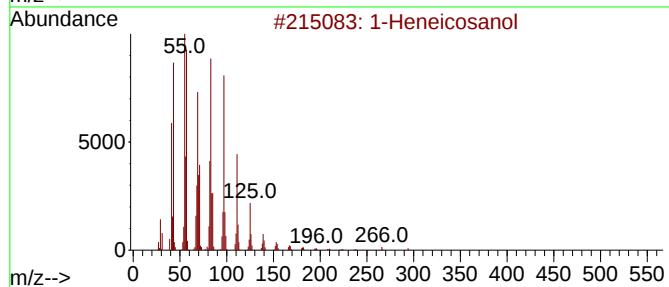
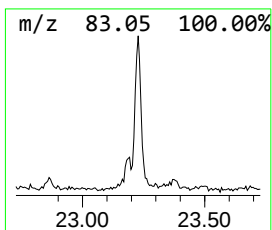
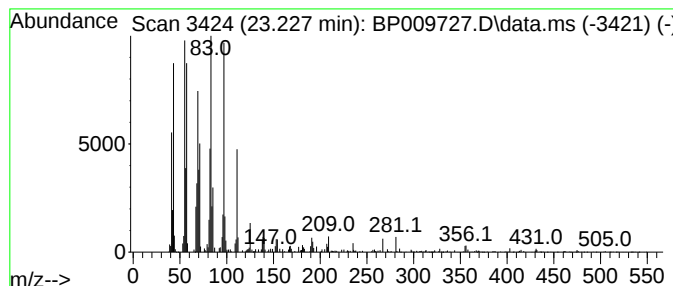
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 1-Heneicosanol Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.227	2.96 ng/ul	323813	Perylene-d12	23.910

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Heneicosanol	312	C21H44O	015594-90-8	91
2			Behenic alcohol	326	C22H46O	000661-19-8	91
3			Cyclotetracosane	336	C24H48	000297-03-0	91
4			Fumaric acid, 2-chloropropyl tri...	374	C20H35ClO4	1010348-57-1	90
5			1-Hexacosene	364	C26H52	018835-33-1	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040822\
Data File : BP009727.D
Acq On : 08 Apr 2022 21:01
Operator : CG/JU
Sample : N2182-08
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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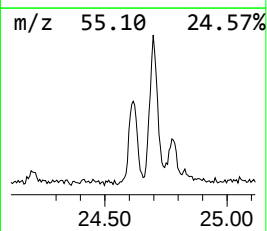
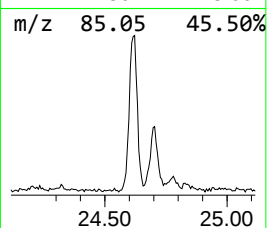
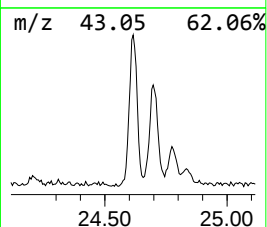
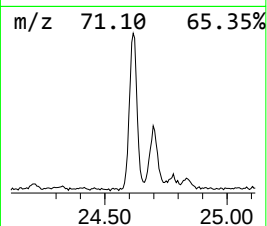
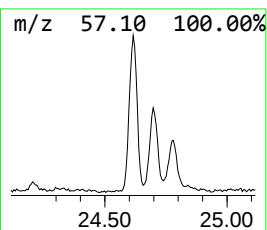
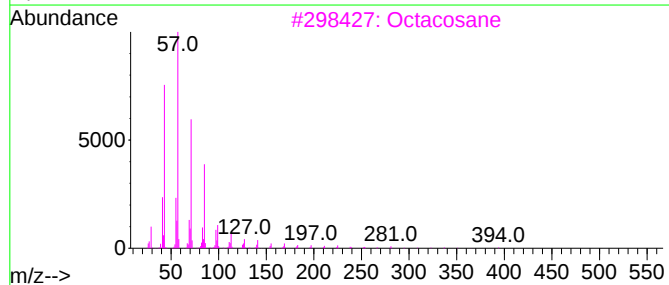
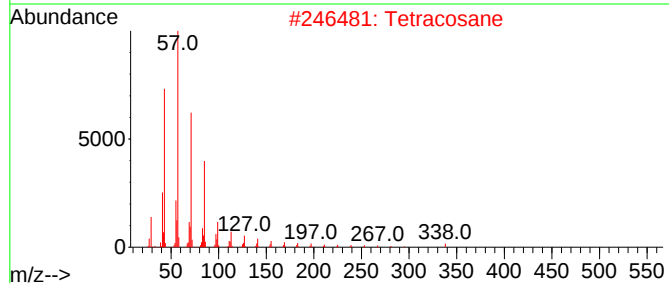
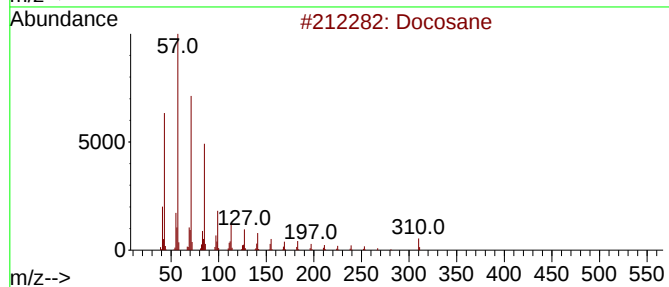
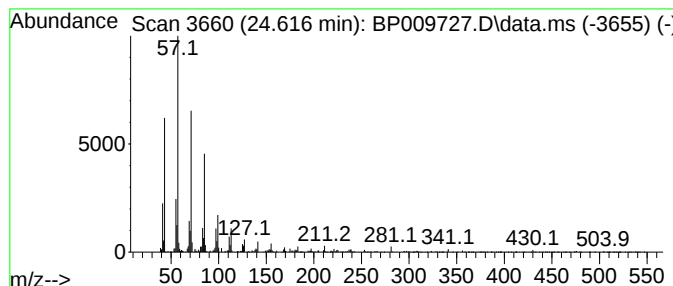
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.616	2.90 ng/ul	316721	Perylene-d12	23.910

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Docosane	310	C22H46	000629-97-0	97
2			Tetracosane	338	C24H50	000646-31-1	90
3			Octacosane	394	C28H58	000630-02-4	90
4			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	90
5			Tetracosane	338	C24H50	000646-31-1	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040822\
Data File : BP009727.D
Acq On : 08 Apr 2022 21:01
Operator : CG/JU
Sample : N2182-08
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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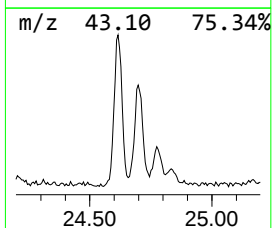
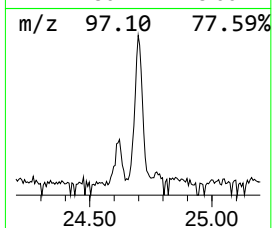
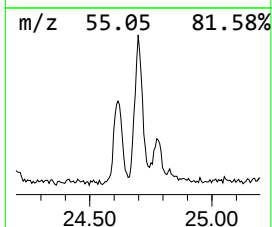
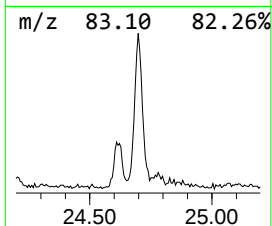
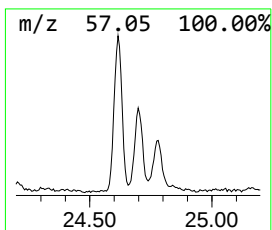
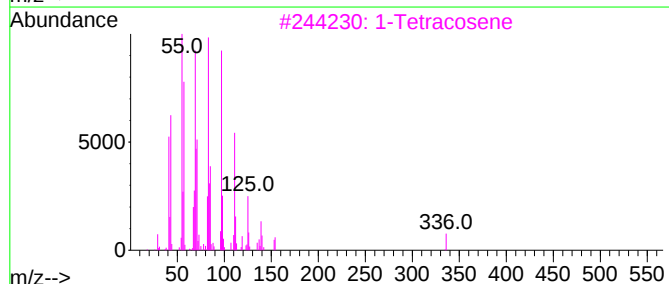
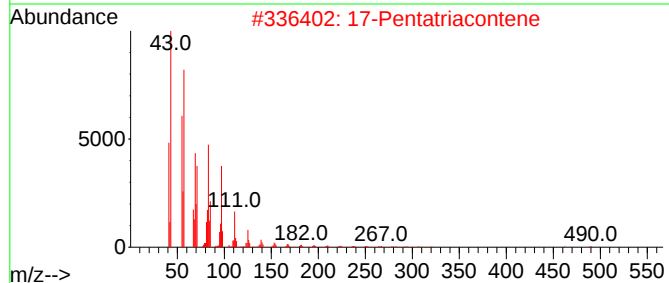
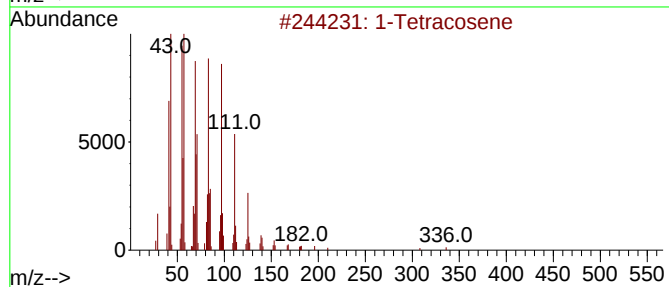
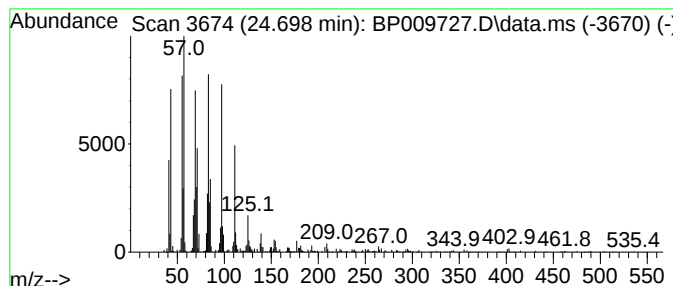
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.698	2.86 ng/ul	313069	Perylene-d12	23.910

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Tetracosene	336	C24H48	010192-32-2	91
2		17-Pentatriacontene	491	C35H70	006971-40-0	86
3		1-Tetracosene	336	C24H48	010192-32-2	83
4		Heptacosyl acetate	438	C29H58O2	1000351-78-2	70
5		Cyclohexane, 1,2,4,5-tetraethyl-...	196	C14H28	061142-24-3	68



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040822\
Data File : BP009727.D
Acq On : 08 Apr 2022 21:01
Operator : CG/JU
Sample : N2182-08
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0R15

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.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
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n-Hexadecanoic ...	18.245	2.5	ng/ul	221999	4	17.357	1797720	20.0
Pentadecafluoro...	21.151	2.5	ng/ul	307650	5	21.439	2430410	20.0
1-Heptacosanol	22.092	2.2	ng/ul	264327	5	21.439	2430410	20.0
1-Heneicosanol	23.227	3.0	ng/ul	323813	6	23.910	2187690	20.0
(DEL) Alkane: S...	24.616	2.9	ng/ul	316721	6	23.910	2187690	20.0
(DEL) Alkane: S...	24.698	2.9	ng/ul	313069	6	23.910	2187690	20.0