

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
 Data File : BP013133.D
 Acq On : 19 Dec 2022 18:55
 Operator : CG/JU
 Sample : N6006-11
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBXZ4

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-BP120722.M
 Title : SVOA CALIBRATION

Signal : TIC: BP013133.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.340	22	26	32	rVB	153377	198897	1.63%	0.106%
2	3.634	73	76	81	rVB	19306	27000	0.22%	0.014%
3	3.769	96	99	113	rBV	762022	1136949	9.30%	0.605%
4	3.875	113	117	124	rVV	96024	137399	1.12%	0.073%
5	3.952	127	130	133	rVV2	15759	22704	0.19%	0.012%
6	3.999	133	138	144	rVB2	84070	129842	1.06%	0.069%
7	4.099	151	155	159	rVB	41310	57376	0.47%	0.031%
8	4.481	216	220	224	rVB3	21073	33415	0.27%	0.018%
9	4.716	256	260	271	rVB4	16560	39156	0.32%	0.021%
10	5.040	309	315	327	rVV	838060	1170970	9.58%	0.623%
11	5.275	352	355	364	rVB2	22651	44603	0.36%	0.024%
12	6.934	631	637	642	rBV6	18152	34748	0.28%	0.018%
13	7.110	658	667	680	rBV	2593938	4688015	38.35%	2.494%
14	7.281	690	696	707	rVB	2187523	3331435	27.26%	1.772%
15	7.481	723	730	740	rBV	2499711	3995987	32.69%	2.125%
16	7.569	742	745	750	rVB4	24880	31586	0.26%	0.017%
17	7.957	803	811	817	rBV	1963824	3122899	25.55%	1.661%
18	8.016	817	821	827	rVB3	19476	34929	0.29%	0.019%
19	8.205	846	853	858	rBV6	21254	42355	0.35%	0.023%
20	8.663	923	931	942	rBV	2551137	4019849	32.89%	2.138%
21	8.787	947	952	958	rVB2	84917	142975	1.17%	0.076%
22	9.122	1002	1009	1019	rVV	2478230	4000086	32.73%	2.128%
23	9.228	1024	1027	1032	rVV4	20031	29572	0.24%	0.016%
24	9.528	1071	1078	1084	rVB	78184	124478	1.02%	0.066%
25	9.851	1123	1133	1144	rBV	2026865	3413571	27.93%	1.816%
26	9.987	1150	1156	1164	rBV3	19129	42327	0.35%	0.023%
27	10.069	1166	1170	1180	rVB9	14602	31446	0.26%	0.017%
28	10.387	1216	1224	1236	rBV	3131809	5174860	42.34%	2.752%
29	10.546	1247	1251	1258	rVB7	11275	24296	0.20%	0.013%
30	10.769	1280	1289	1297	rBV	2742467	4584977	37.51%	2.439%
31	10.910	1302	1313	1334	rVV	1948804	3555087	29.09%	1.891%
32	11.298	1374	1379	1387	rVB8	21650	41880	0.34%	0.022%
33	11.781	1455	1461	1466	rBV7	15073	30319	0.25%	0.016%
34	12.081	1507	1512	1517	rVV2	18651	28169	0.23%	0.015%
35	12.181	1523	1529	1538	rVB	69685	106474	0.87%	0.057%
36	12.293	1544	1548	1553	rVV	34255	53367	0.44%	0.028%

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Smoothing : OFF

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
 Title : SVOA CALIBRATION

37	12.351	1553	1558	1564	rVB	57178	87577	0.72%	0.047%
38	12.457	1568	1576	1591	rBV	1030839	1687781	13.81%	0.898%
39	12.922	1650	1655	1659	rBV7	15118	24943	0.20%	0.013%
40	12.987	1662	1666	1669	rBV4	16028	23098	0.19%	0.012%
41	13.128	1686	1690	1694	rVV	20820	27146	0.22%	0.014%
42	13.410	1728	1738	1745	rVB	118404	177284	1.45%	0.094%
43	13.487	1747	1751	1756	rBV5	14047	25596	0.21%	0.014%
44	13.587	1761	1768	1773	rVB3	41709	74793	0.61%	0.040%
45	13.922	1819	1825	1830	rBV9	27710	51653	0.42%	0.027%
46	14.004	1831	1839	1847	rBV	5913847	7776870	63.63%	4.136%
47	14.069	1848	1850	1854	rVB	28533	30403	0.25%	0.016%
48	14.122	1854	1859	1864	rVB5	40367	61078	0.50%	0.032%
49	14.187	1864	1870	1874	rBV	86849	121808	1.00%	0.065%
50	14.228	1874	1877	1881	rBV3	19750	29515	0.24%	0.016%
51	14.292	1881	1888	1899	rVV	6463536	9235277	75.56%	4.912%
52	14.487	1913	1921	1925	rVB	78984	113352	0.93%	0.060%
53	14.598	1931	1940	1947	rBV	4288695	6079721	49.74%	3.234%
54	14.792	1966	1973	1993	rBV	1909087	3080299	25.20%	1.638%
55	14.945	1995	1999	2004	rVV3	54952	95248	0.78%	0.051%
56	15.010	2004	2010	2021	rVB6	82650	173549	1.42%	0.092%
57	15.387	2069	2074	2078	rBV4	17547	36979	0.30%	0.020%
58	15.434	2080	2082	2087	rVB2	22727	30597	0.25%	0.016%
59	15.592	2102	2109	2115	rBV	8055558	11542442	94.43%	6.139%
60	15.645	2115	2118	2123	rVB2	48146	67384	0.55%	0.036%
61	15.710	2123	2129	2140	rBV	2405374	3204219	26.21%	1.704%
62	15.834	2146	2150	2156	rVB2	40464	68068	0.56%	0.036%
63	15.957	2165	2171	2183	rBV	5191367	6790948	55.56%	3.612%
64	16.298	2226	2229	2232	rBV5	16901	24953	0.20%	0.013%
65	16.328	2232	2234	2239	rVB6	24338	25352	0.21%	0.013%
66	16.439	2249	2253	2258	rBV4	32465	50315	0.41%	0.027%
67	16.481	2258	2260	2263	rVV2	16504	23379	0.19%	0.012%
68	16.522	2263	2267	2272	rVB	313792	408220	3.34%	0.217%
69	16.733	2300	2303	2311	rVB4	36644	55334	0.45%	0.029%
70	16.886	2325	2329	2333	rBV2	33619	47357	0.39%	0.025%
71	17.004	2345	2349	2351	rBV5	17355	24084	0.20%	0.013%
72	17.233	2384	2388	2396	rBV3	71703	120594	0.99%	0.064%
73	17.304	2396	2400	2402	rVV3	39199	51721	0.42%	0.028%
74	17.351	2402	2408	2413	rVV	4653355	6733151	55.09%	3.581%
75	17.392	2413	2415	2419	rVV	191363	251469	2.06%	0.134%
76	17.451	2419	2425	2437	rVV	7396401	10948916	89.58%	5.824%

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 Title : SVOA CALIBRATION

77	17.545	2437	2441	2449	rVB4	97005	155637	1.27%	0.083%
78	18.010	2517	2520	2523	rBV4	28833	32133	0.26%	0.017%
79	18.063	2525	2529	2531	rBV5	27167	41846	0.34%	0.022%
80	18.110	2531	2537	2542	rBV4	92045	179407	1.47%	0.095%
81	18.163	2543	2546	2551	rBV	156664	195736	1.60%	0.104%
82	18.228	2552	2557	2566	rBV	1455961	2110951	17.27%	1.123%
83	18.510	2602	2605	2610	rVB5	56572	78531	0.64%	0.042%
84	18.698	2635	2637	2641	rVB3	48311	57907	0.47%	0.031%
85	19.110	2702	2707	2715	rBV6	101216	214630	1.76%	0.114%
86	19.210	2721	2724	2726	rBV2	34306	47928	0.39%	0.025%
87	19.263	2729	2733	2738	rVB2	123279	182338	1.49%	0.097%
88	19.357	2739	2749	2754	rBV2	467154	1046327	8.56%	0.557%
89	19.451	2761	2765	2780	rBV	453507	742599	6.08%	0.395%
90	19.686	2799	2805	2817	rBV	8656224	12222928	100.00%	6.501%
91	20.192	2888	2891	2896	rVB2	139384	162922	1.33%	0.087%
92	21.127	3046	3050	3056	rBV	1164283	1942205	15.89%	1.033%
93	21.439	3097	3103	3108	rBV	4329063	5902968	48.29%	3.140%
94	23.139	3386	3392	3398	rBV	4259421	7143987	58.45%	3.800%
95	23.757	3491	3497	3512	rVB2	5040507	10847855	88.75%	5.770%
96	23.915	3518	3524	3545	rVB2	2772130	6098050	49.89%	3.243%
97	24.433	3605	3612	3622	rVB2	2876434	6078941	49.73%	3.233%
98	24.539	3624	3630	3641	rVB	3212485	7093183	58.03%	3.773%
99	26.457	3950	3956	3975	rVB2	1056648	3432344	28.08%	1.826%
100	28.509	4298	4305	4325	rVB3	2369044	8831221	72.25%	4.697%

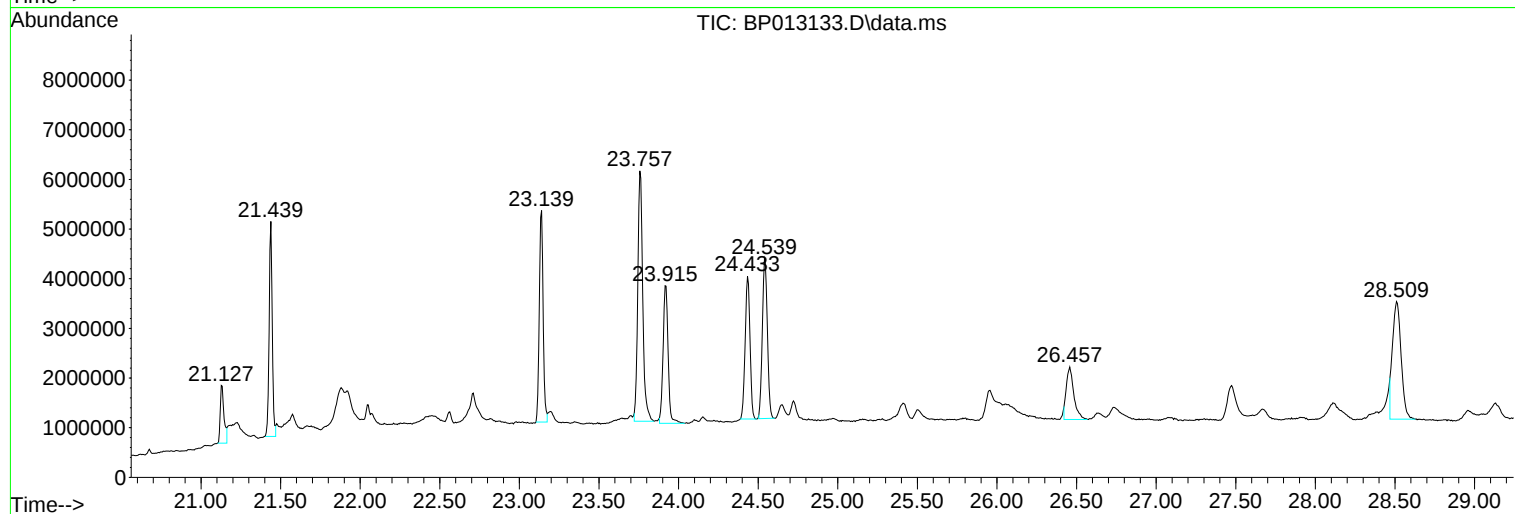
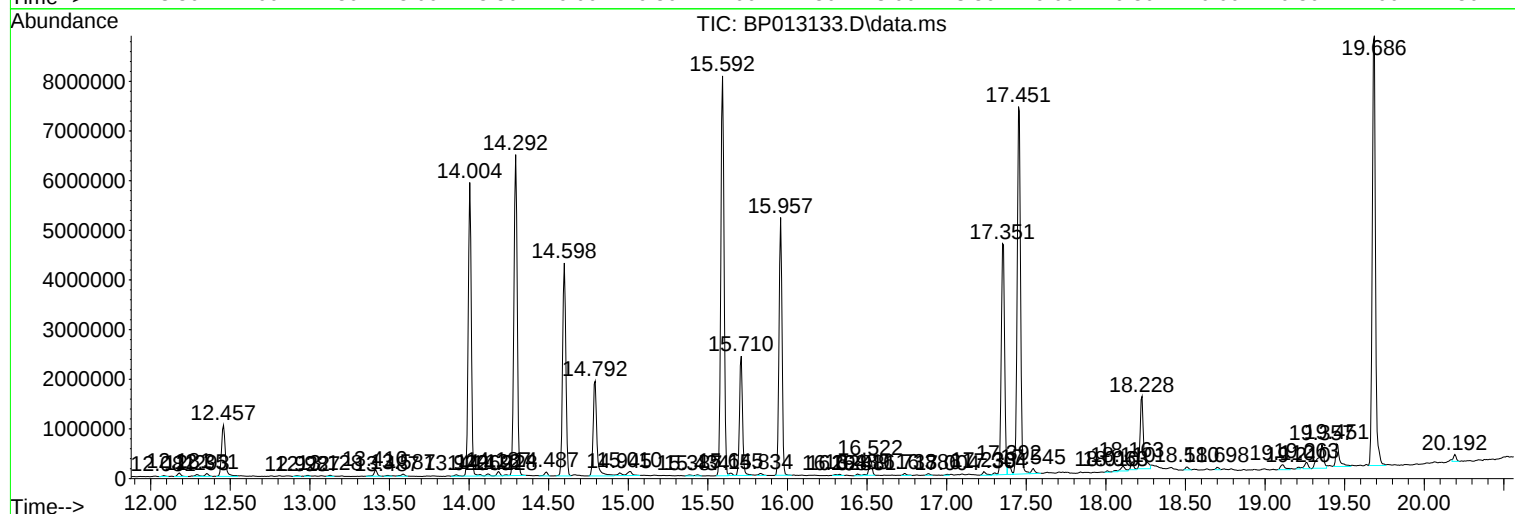
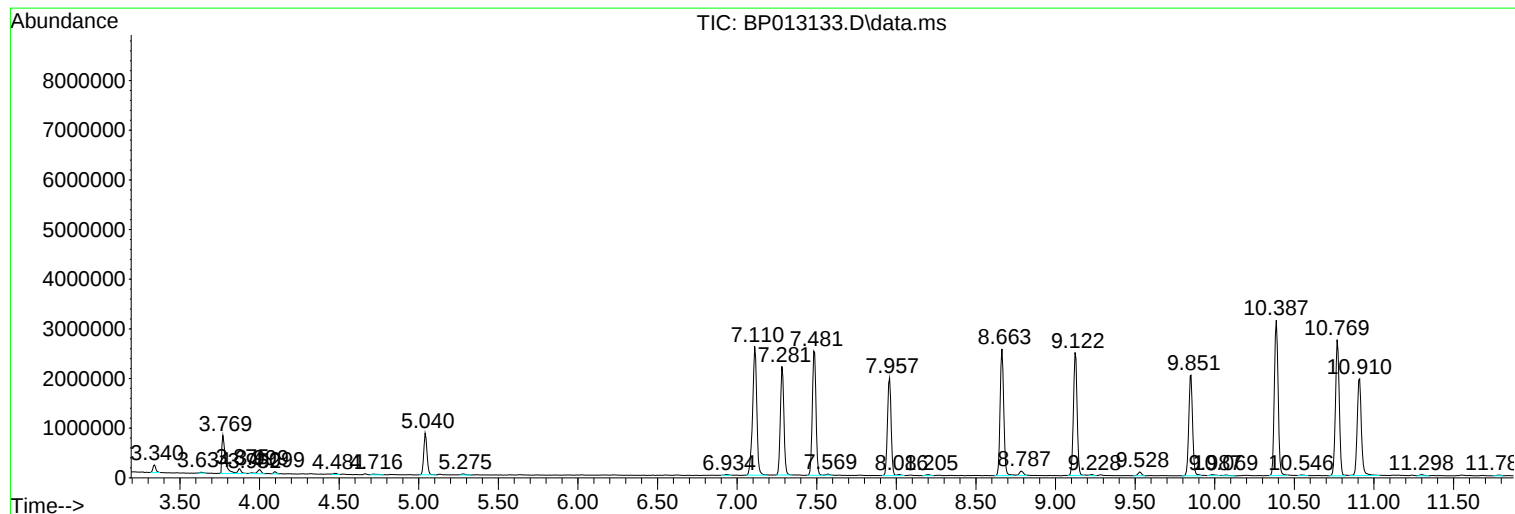
Sum of corrected areas: 188009045

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BNA_P
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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



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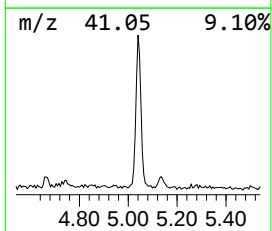
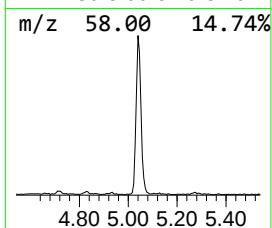
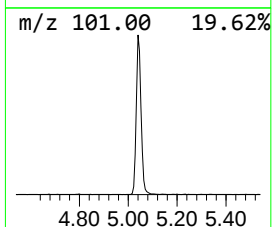
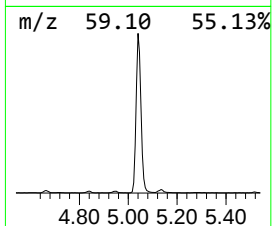
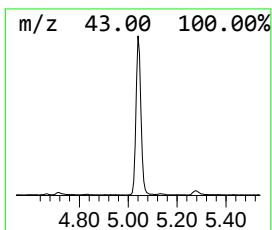
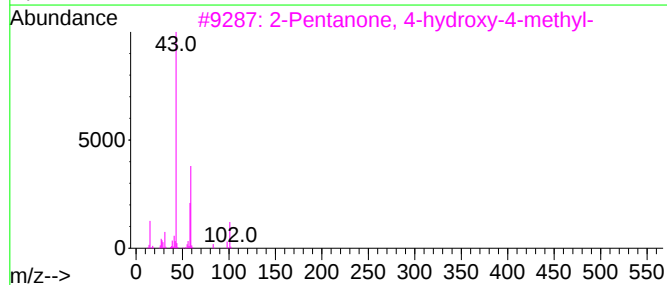
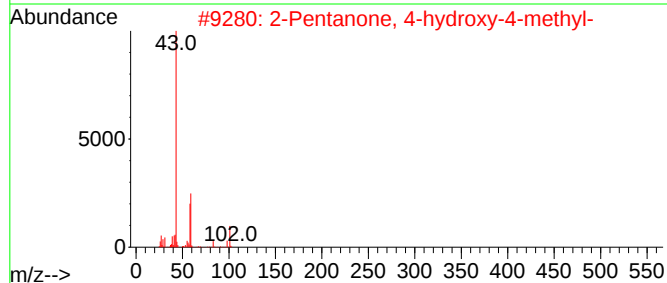
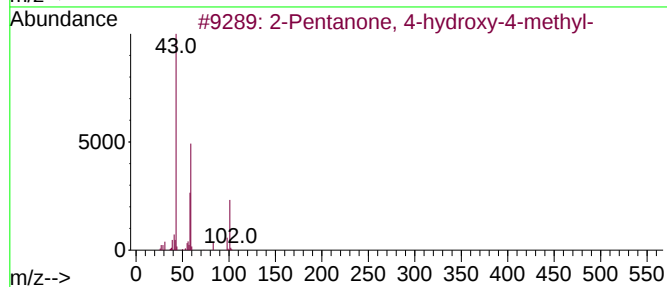
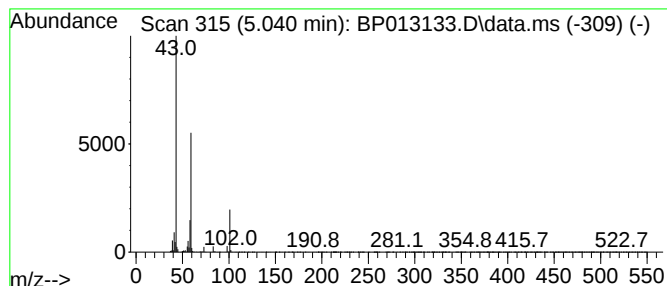
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.040	7.50 ng/ul	1170970	1,4-Dichlorobenzene-d4	7.957

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56		
2	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50		
3	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	40		
4	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	40		
5	Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	32		



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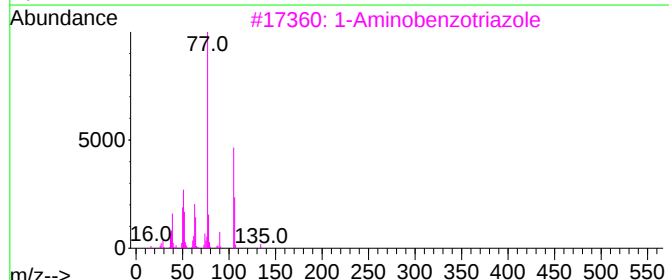
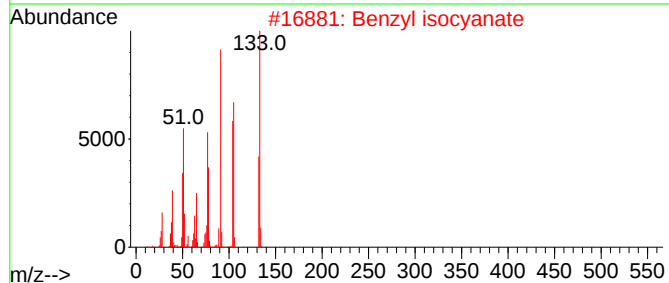
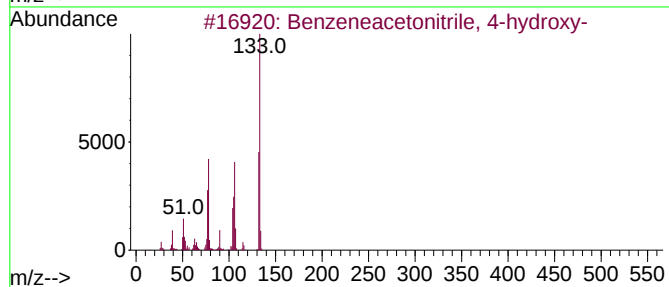
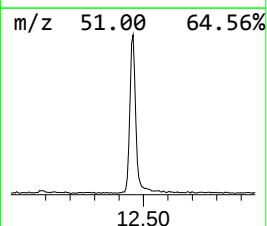
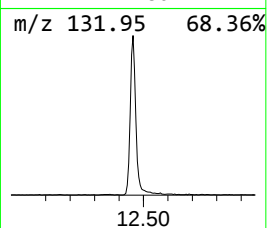
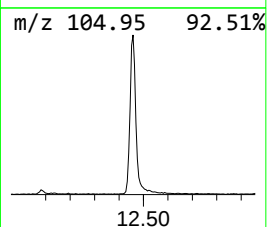
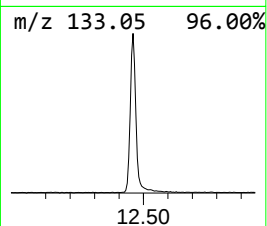
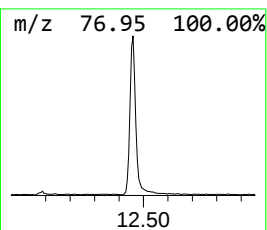
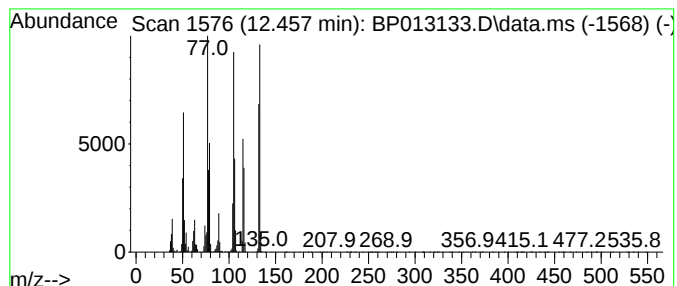
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Quant Title : SVOA CALIBRATION

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

Peak Number 2 unknown-01 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.457	7.36 ng/ul	1687780	Naphthalene-d8	10.769

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzeneacetonitrile, 4-hydroxy-	133	C8H7NO	014191-95-8	41
2			Benzyl isocyanate	133	C8H7NO	003173-56-6	30
3			1-Aminobenzotriazole	134	C6H6N4	001614-12-6	27
4			2-Propen-1-one, 1-phenyl-	132	C9H8O	000768-03-6	27
5			5-Aza-2H-indazole, N-methyl	133	C7H7N3	1010508-64-4	25



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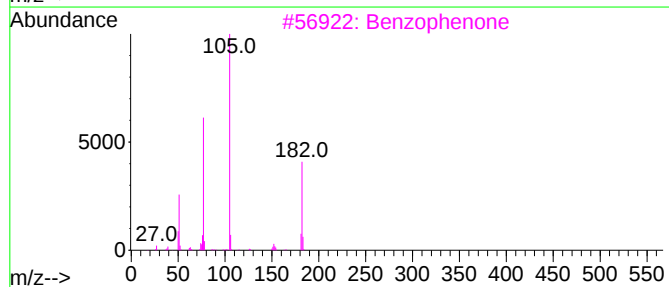
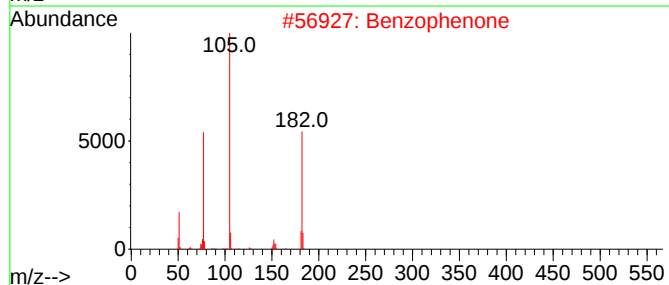
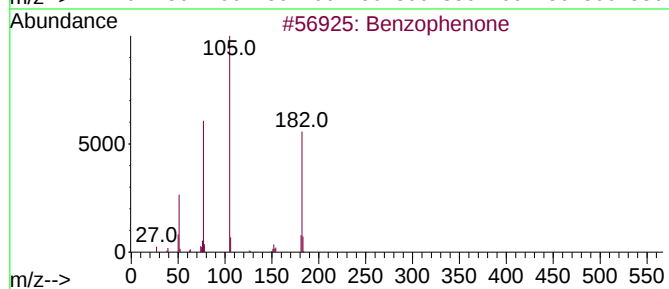
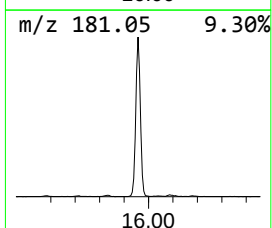
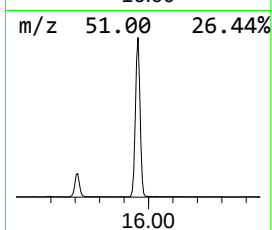
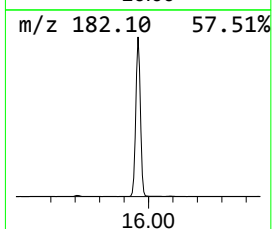
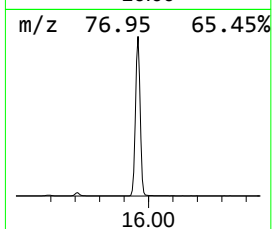
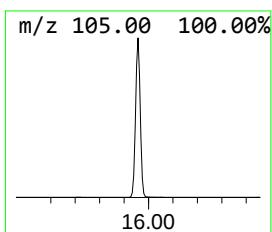
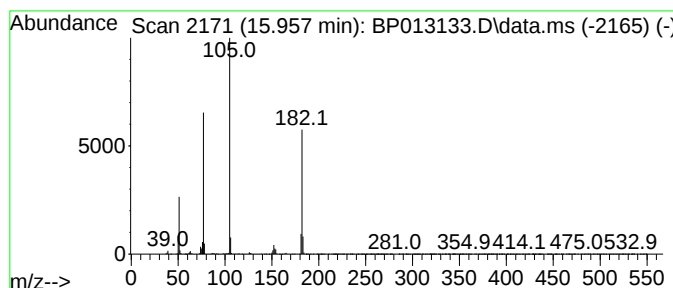
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Benzophenone Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.957	22.34 ng/ul	6790950	Acenaphthene-d10	14.598

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	93
5			Benzophenone	182	C13H10O	000119-61-9	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013133.D
Acq On : 19 Dec 2022 18:55
Operator : CG/JU
Sample : N6006-11
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBXZ4

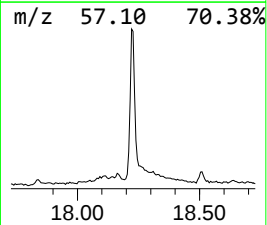
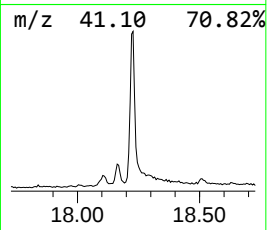
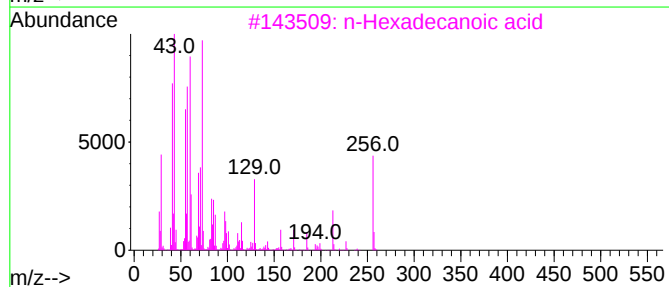
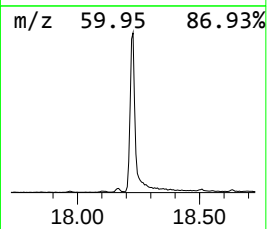
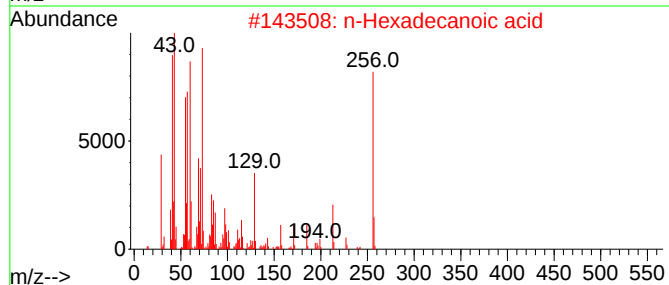
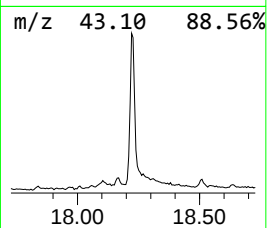
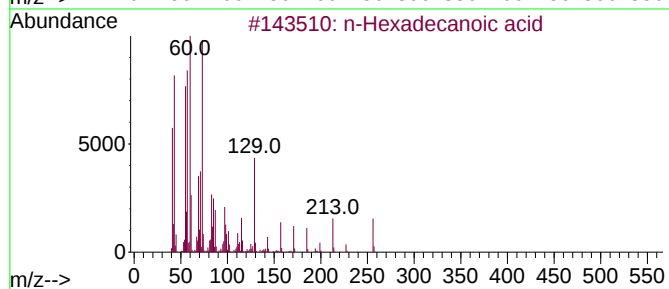
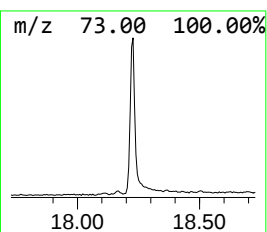
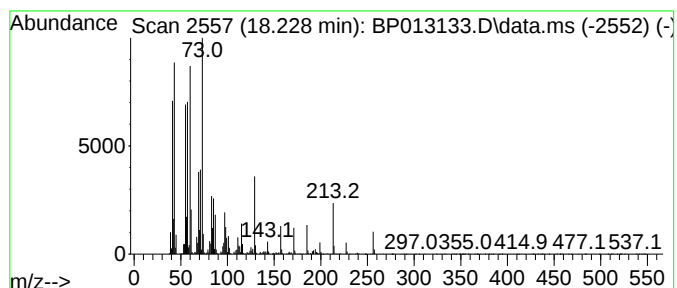
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 n-Hexadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.227	6.27 ng/ul	2110950	Phenanthrene-d10	17.351

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	97
5			Tridecanoic acid	214	C13H26O2	000638-53-9	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013133.D
Acq On : 19 Dec 2022 18:55
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Sample : N6006-11
Misc :
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Instrument :
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ClientSampleId :
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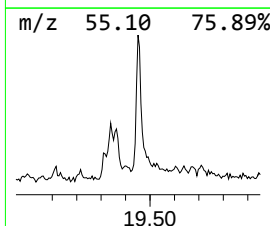
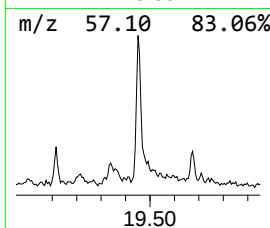
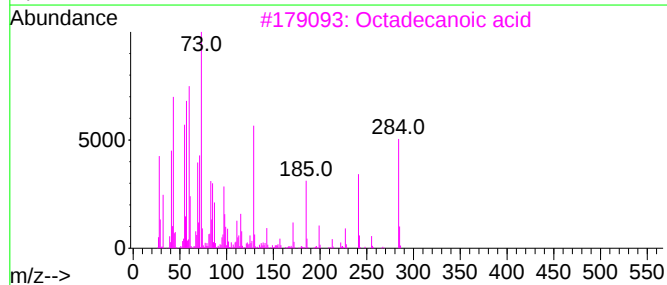
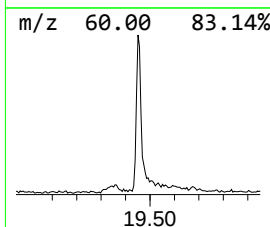
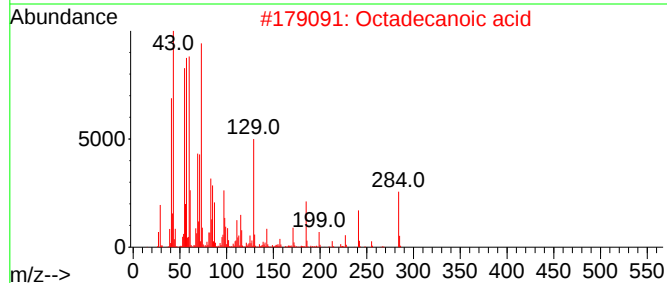
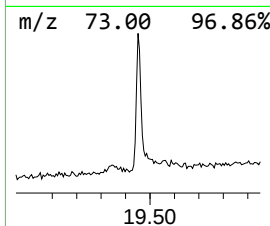
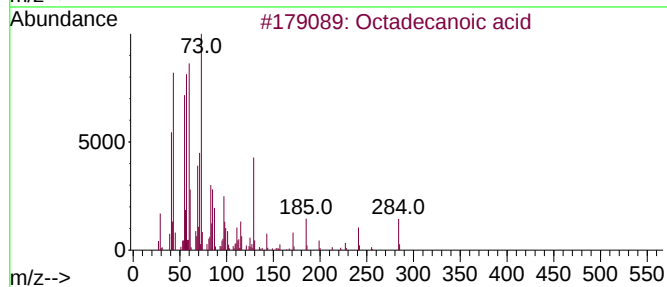
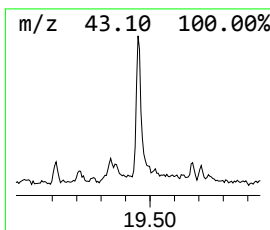
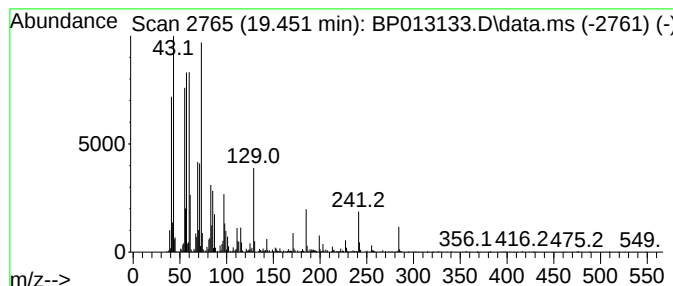
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Octadecanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.451	2.52 ng/ul	742599	Chrysene-d12	21.439

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	99
3			Octadecanoic acid	284	C18H36O2	000057-11-4	99
4			Octadecanoic acid	284	C18H36O2	000057-11-4	95
5			Octadecanoic acid	284	C18H36O2	000057-11-4	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013133.D
Acq On : 19 Dec 2022 18:55
Operator : CG/JU
Sample : N6006-11
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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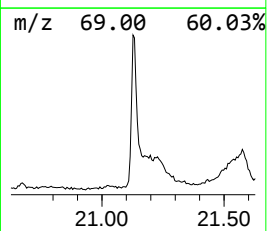
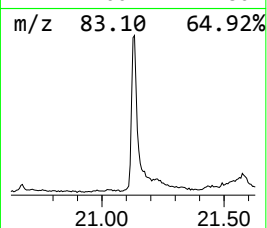
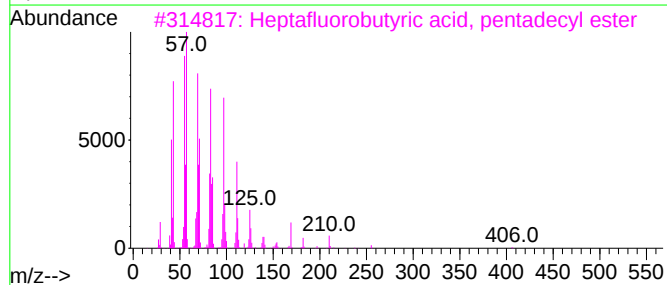
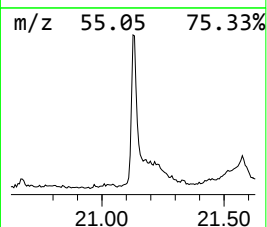
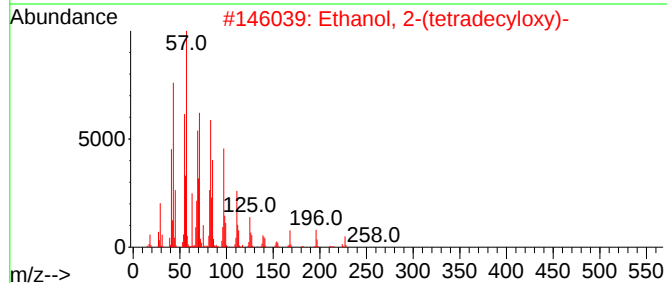
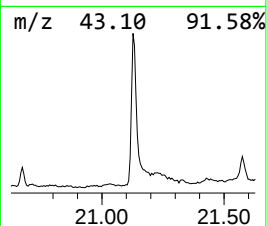
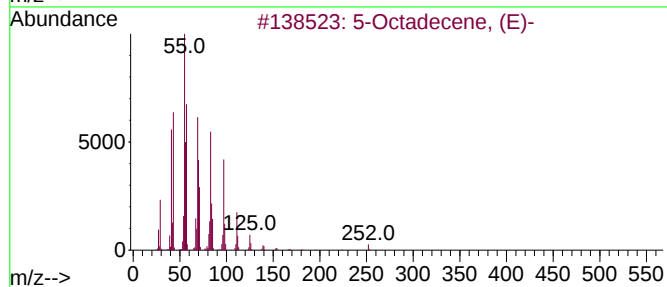
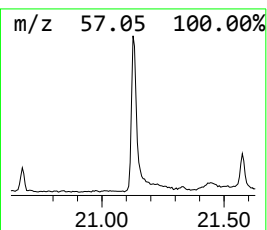
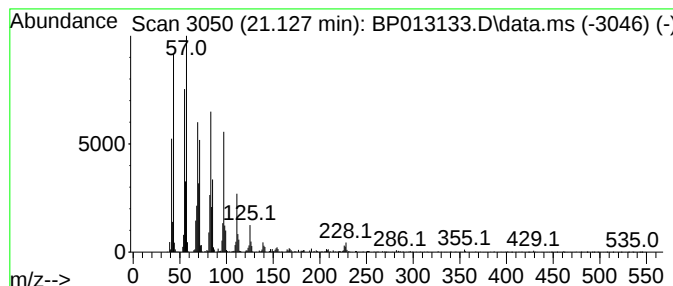
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.127	6.58 ng/ul	1942210	Chrysene-d12	21.439

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			5-Octadecene, (E)-	252	C18H36	007206-21-5	95
2			Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	95
3			Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	94
4			Heptafluorobutyric acid, hexadec...	438	C20H33F7O2	006385-15-5	94
5			Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013133.D
Acq On : 19 Dec 2022 18:55
Operator : CG/JU
Sample : N6006-11
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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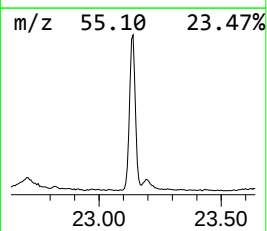
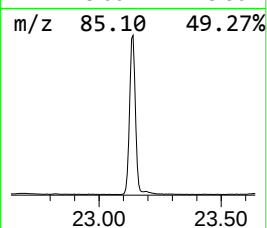
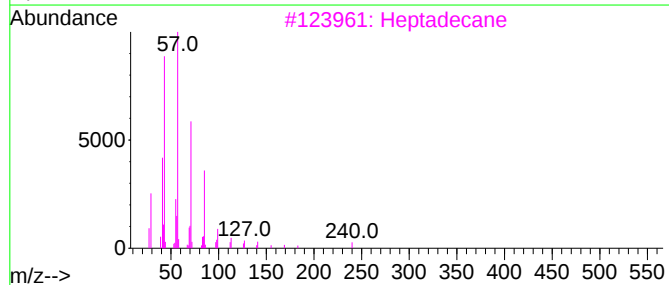
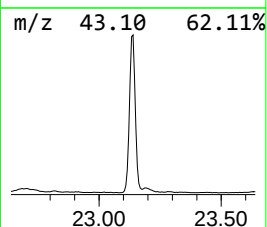
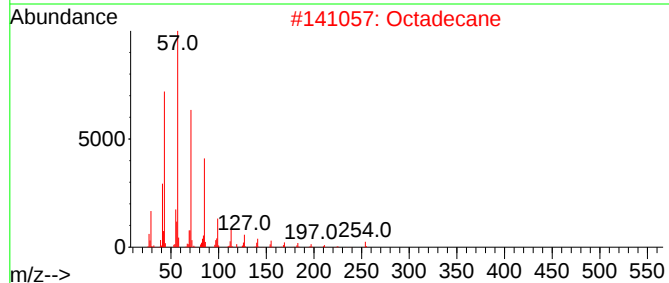
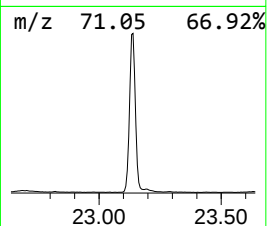
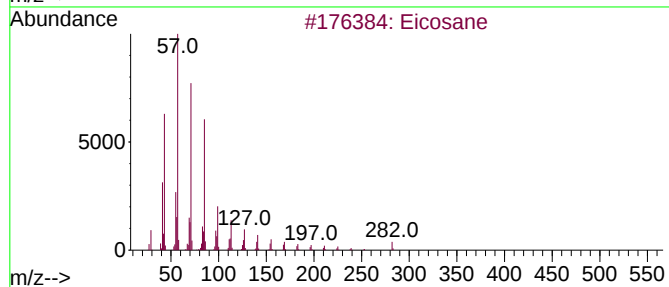
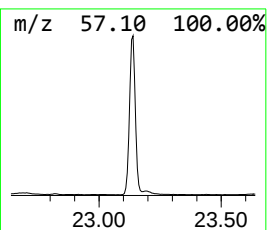
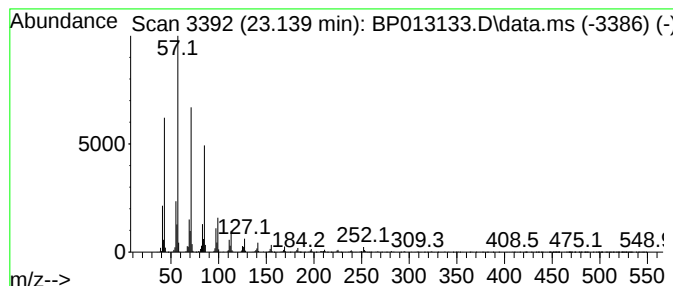
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.139	23.43 ng/ul	7143990	Perylene-d12	23.921

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosane	282	C20H42	000112-95-8	96
2			Octadecane	254	C18H38	000593-45-3	95
3			Heptadecane	240	C17H36	000629-78-7	94
4			2-Bromo dodecane	248	C12H25Br	013187-99-0	93
5			Tetratetracontane	619	C44H90	007098-22-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013133.D
Acq On : 19 Dec 2022 18:55
Operator : CG/JU
Sample : N6006-11
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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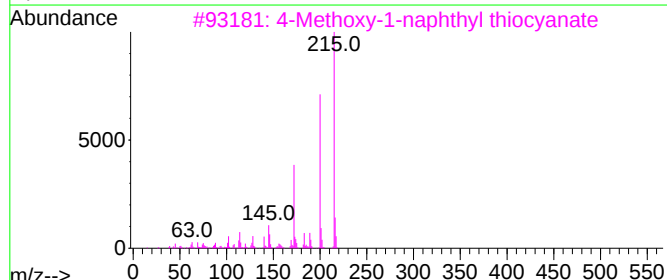
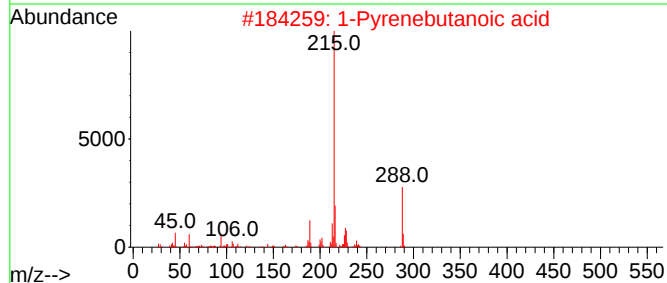
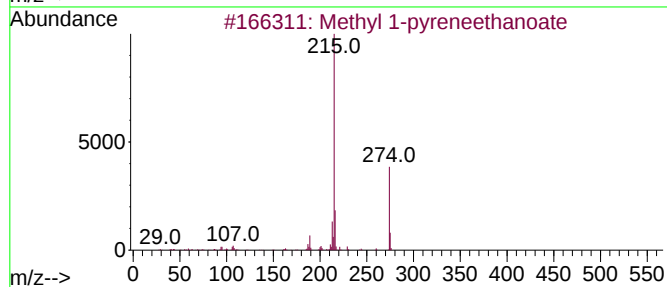
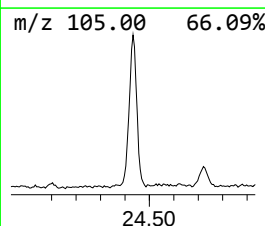
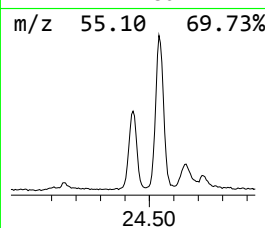
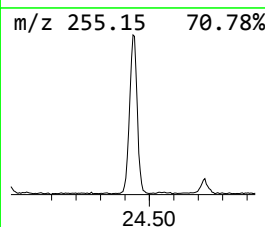
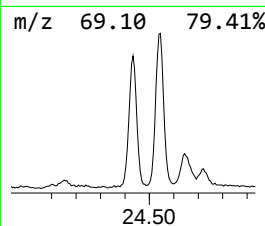
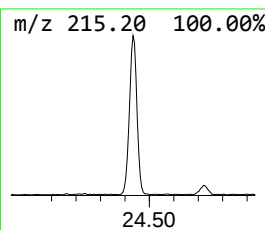
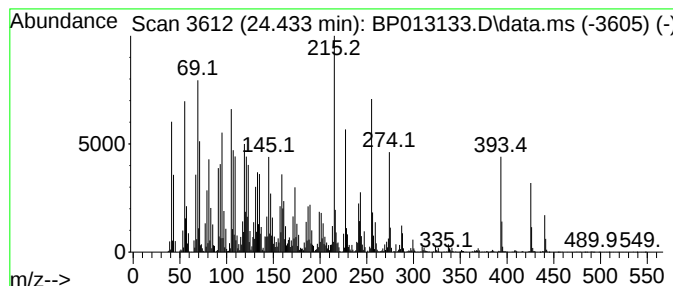
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 unknown-02 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.433	19.94 ng/ul	6078940	Perylene-d12	23.921

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methyl 1-pyreneethanoate	274	C19H14O2	073654-19-0	42
2			1-Pyrenebutanoic acid	288	C20H16O2	003443-45-6	20
3			4-Methoxy-1-naphthyl thiocyanate	215	C12H9NOS	066231-77-4	20
4			Indanidine	215	C11H13N5	085392-79-6	20
5			2(3H)-Benzofuranone, 3-[3-(dimet...	215	C13H13NO2	056771-83-6	18



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013133.D
Acq On : 19 Dec 2022 18:55
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Sample : N6006-11
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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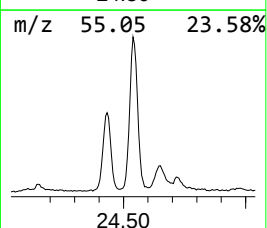
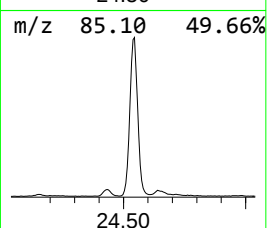
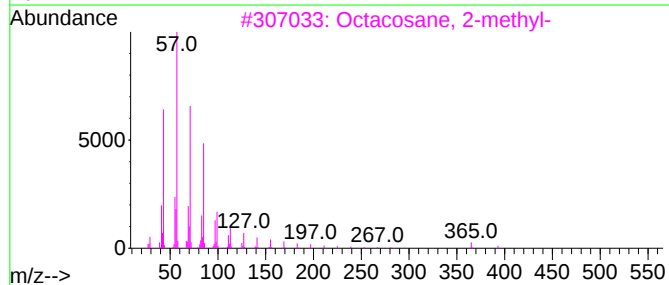
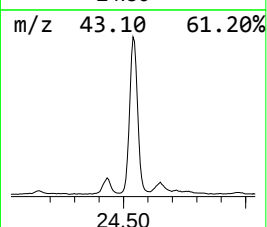
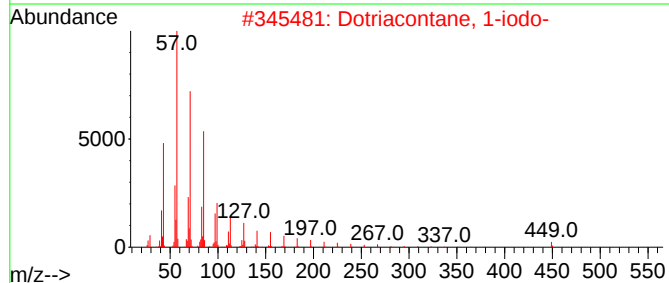
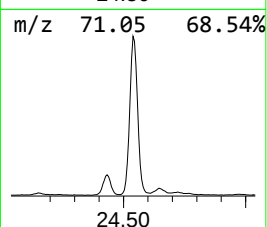
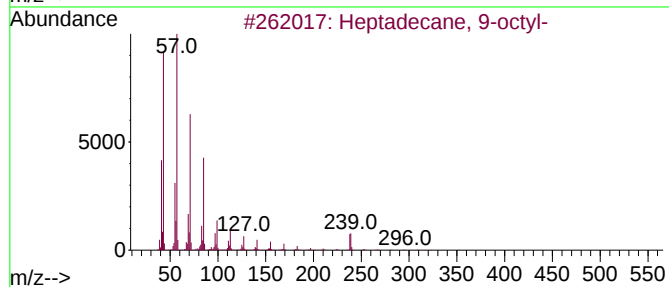
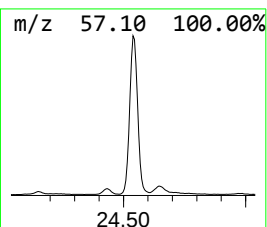
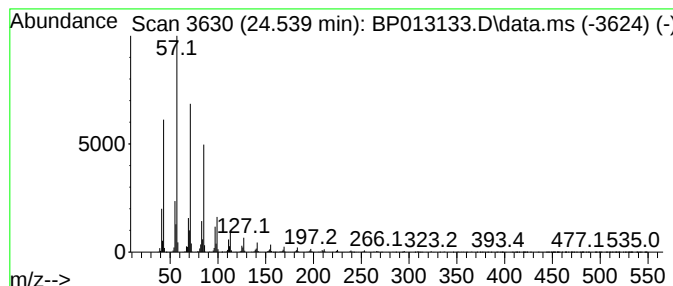
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.539	23.26 ng/ul	7093180	Perylene-d12	23.921

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	95
2		Dotriacontane, 1-iodo-	576	C32H65I	1000406-32-4	91
3		Octacosane, 2-methyl-	408	C29H60	001560-98-1	91
4		Tetratetracontane	619	C44H90	007098-22-8	91
5		Heneicosane	296	C21H44	000629-94-7	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013133.D
Acq On : 19 Dec 2022 18:55
Operator : CG/JU
Sample : N6006-11
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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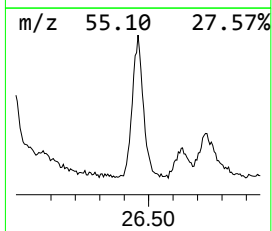
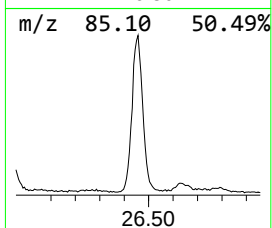
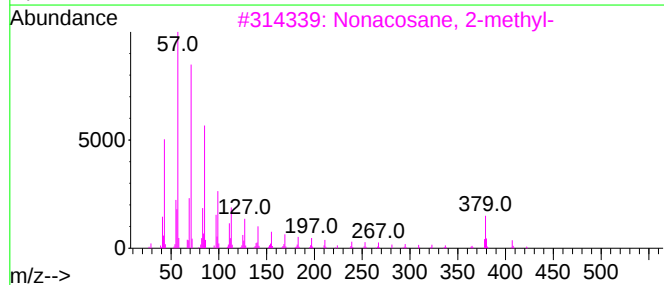
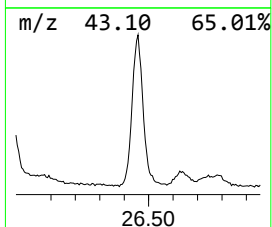
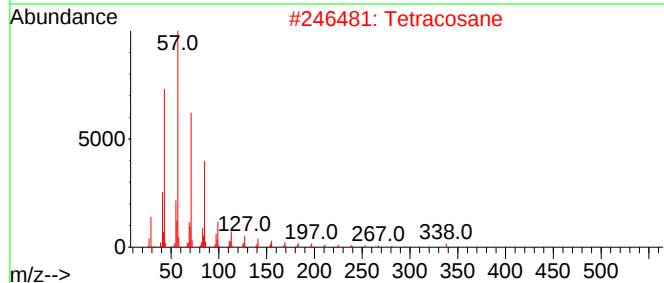
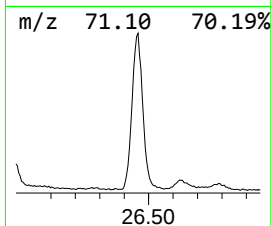
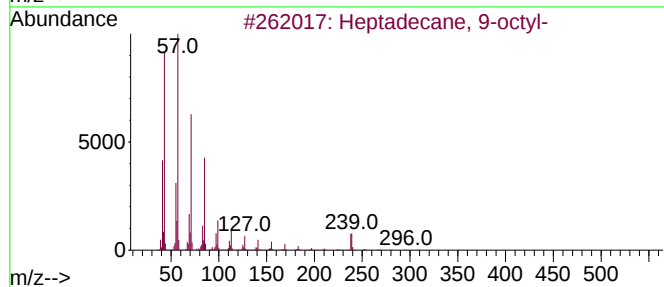
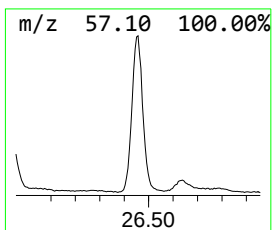
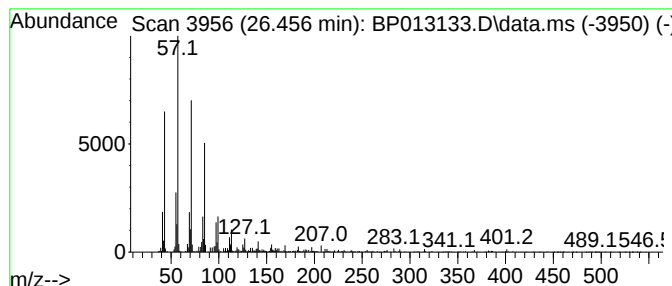
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.456	11.26 ng/ul	3432340	Perylene-d12	23.921

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	93
2			Tetracosane	338	C24H50	000646-31-1	93
3			Nonacosane, 2-methyl-	422	C30H62	001560-75-4	93
4			Tetracosane	338	C24H50	000646-31-1	93
5			Eicosane, 10-methyl-	296	C21H44	054833-23-7	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013133.D
Acq On : 19 Dec 2022 18:55
Operator : CG/JU
Sample : N6006-11
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBXZ4

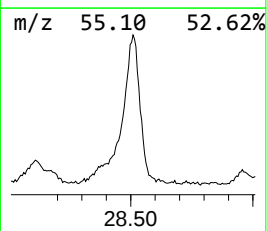
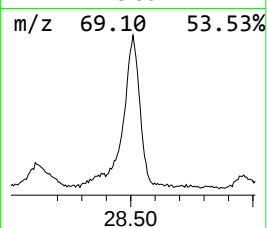
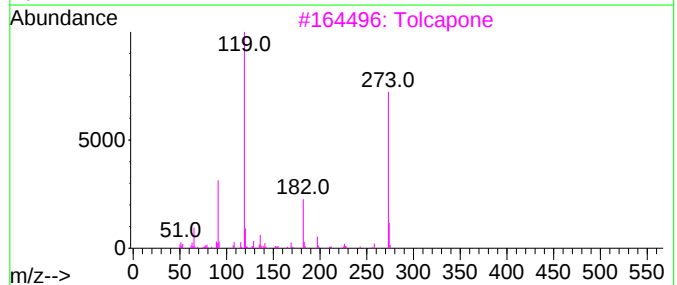
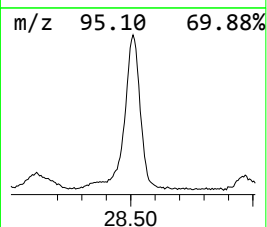
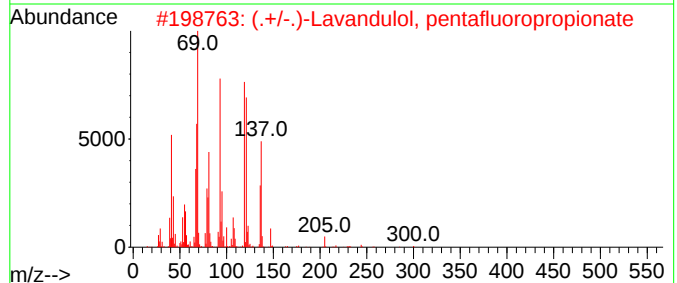
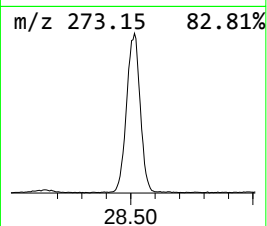
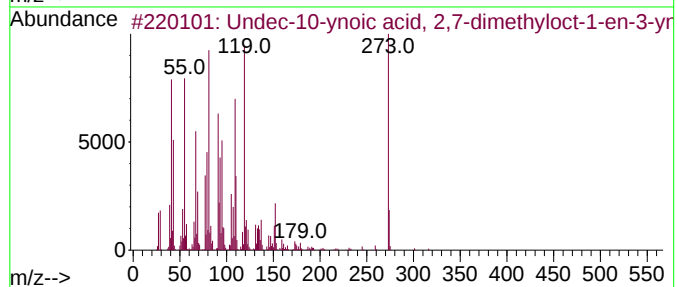
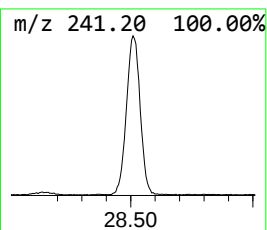
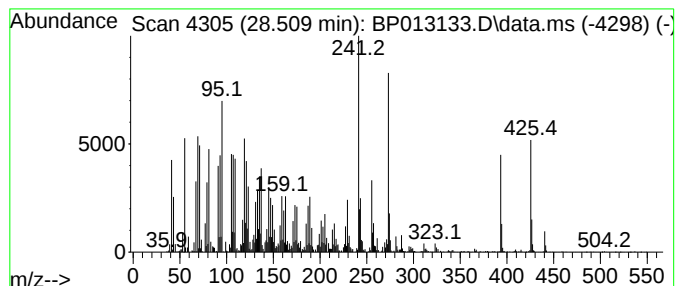
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 unknown-03 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
28.509	28.96 ng/ul	8831220	Perylene-d12	23.921

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Undec-10-ynoic acid, 2,7-dimethy...	316	C21H32O2	1000406-96-7	35
2			(.+/-)-Lavandulol, pentafluorop...	300	C13H17F5O2	1000352-63-1	25
3			Tolcapone	273	C14H11NO5	134308-13-7	25
4			11-Oxo-.alpha.-amyrin	440	C30H48O2	002118-90-3	25
5			6-(Trifluoromethoxy)-1H-indole-2...	273	C12H10F3NO3	1000499-75-7	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
 Data File : BP013133.D
 Acq On : 19 Dec 2022 18:55
 Operator : CG/JU
 Sample : N6006-11
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBXZ4

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.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
2-Pentanone, 4-...	5.040	7.5	ng/u1	1170970	1	7.957	3122900	20.0
unknown-01	12.457	7.4	ng/u1	1687780	2	10.769	4584980	20.0
Benzophenone	15.957	22.3	ng/u1	6790950	3	14.598	6079720	20.0
n-Hexadecanoic ...	18.227	6.3	ng/u1	2110950	4	17.351	6733150	20.0
Octadecanoic acid	19.451	2.5	ng/u1	742599	5	21.439	5902970	20.0
(DEL) Alkane: S...	21.127	6.6	ng/u1	1942210	5	21.439	5902970	20.0
(DEL) Alkane: S...	23.139	23.4	ng/u1	7143990	6	23.921	6098050	20.0
unknown-02	24.433	19.9	ng/u1	6078940	6	23.921	6098050	20.0
(DEL) Alkane: S...	24.539	23.3	ng/u1	7093180	6	23.921	6098050	20.0
(DEL) Alkane: S...	26.456	11.3	ng/u1	3432340	6	23.921	6098050	20.0
unknown-03	28.509	29.0	ng/u1	8831220	6	23.921	6098050	20.0