

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111022\
 Data File : BP012604.D
 Acq On : 10 Nov 2022 17:50
 Operator : CG/JU
 Sample : N5493-06
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 H0AC5

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-BP110922.M
 Title : SVOA CALIBRATION

Signal : TIC: BP012604.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.222	37	40	45	rBV	46599	63024	0.74%	0.080%
2	3.652	111	113	133	rBV	157856	326429	3.82%	0.415%
3	3.852	143	147	155	rVB2	20355	36033	0.42%	0.046%
4	3.952	160	164	172	rVB	21767	32324	0.38%	0.041%
5	4.334	225	229	234	rVB8	4679	8618	0.10%	0.011%
6	6.952	667	674	685	rBV	202945	379059	4.43%	0.482%
7	7.099	691	699	716	rBV	844075	1445074	16.89%	1.839%
8	7.299	726	733	743	rBV	777614	1318791	15.42%	1.678%
9	7.757	804	811	822	rBV	971561	1666055	19.47%	2.120%
10	8.487	928	935	958	rBV	628542	1159271	13.55%	1.475%
11	8.928	1003	1010	1030	rBV	979654	1749473	20.45%	2.227%
12	9.063	1030	1033	1045	rVB	5825	18708	0.22%	0.024%
13	9.304	1067	1074	1080	rVB2	26599	40573	0.47%	0.052%
14	9.651	1125	1133	1145	rBV	785033	1363728	15.94%	1.736%
15	10.193	1216	1225	1245	rBV	1040097	2021189	23.63%	2.572%
16	10.557	1279	1287	1302	rBV	1569643	2699473	31.55%	3.436%
17	10.716	1306	1314	1339	rBV	782011	1605467	18.77%	2.043%
18	11.993	1525	1531	1537	rBV7	8870	16700	0.20%	0.021%
19	12.163	1555	1560	1567	rVB3	16816	32740	0.38%	0.042%
20	12.416	1598	1603	1610	rBV2	12312	18186	0.21%	0.023%
21	13.228	1736	1741	1744	rBV5	6991	10180	0.12%	0.013%
22	13.304	1748	1754	1760	rVV2	28326	46406	0.54%	0.059%
23	13.381	1764	1767	1776	rVB8	6749	13990	0.16%	0.018%
24	13.834	1835	1844	1858	rBV	2458809	3733649	43.64%	4.752%
25	13.934	1859	1861	1867	rVB5	21021	33003	0.39%	0.042%
26	14.104	1883	1890	1914	rBV	2709849	4293559	50.19%	5.464%
27	14.298	1921	1923	1928	rVB6	9970	13123	0.15%	0.017%
28	14.416	1936	1943	1955	rBV	2598923	3976781	46.48%	5.061%
29	14.687	1983	1989	1996	rBV2	91800	235748	2.76%	0.300%
30	14.839	2012	2015	2022	rVB7	18605	29982	0.35%	0.038%
31	15.075	2051	2055	2061	rVB3	22331	30659	0.36%	0.039%
32	15.210	2075	2078	2081	rBV2	7547	8854	0.10%	0.011%
33	15.257	2081	2086	2090	rVB5	9420	14187	0.17%	0.018%
34	15.416	2106	2113	2130	rBV	3559559	5731883	67.00%	7.295%
35	15.551	2130	2136	2150	rVV	1216871	1872826	21.89%	2.383%
36	15.651	2150	2153	2159	rVB4	24649	42110	0.49%	0.054%

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37	16.145	2231	2237	2243	rBV4	10870	25680	0.30%	0.033%
38	16.581	2307	2311	2319	rBV2	38187	64043	0.75%	0.082%
39	16.939	2368	2372	2376	rBV4	14923	25349	0.30%	0.032%
40	17.181	2406	2413	2423	rBV	3168792	4742979	55.44%	6.036%
41	17.281	2423	2430	2440	rBV	4524517	6848232	80.05%	8.716%
42	17.469	2459	2462	2468	rVB4	14081	19444	0.23%	0.025%
43	17.604	2478	2485	2493	rBV4	27553	71892	0.84%	0.091%
44	17.845	2522	2526	2532	rBV	208435	261513	3.06%	0.333%
45	17.951	2540	2544	2552	rVB8	18812	33500	0.39%	0.043%
46	18.069	2559	2564	2569	rBV	432383	648206	7.58%	0.825%
47	18.951	2712	2714	2715	rBV2	19038	15385	0.18%	0.020%
48	18.980	2715	2719	2730	rVB	434311	537510	6.28%	0.684%
49	19.110	2735	2741	2746	rBV2	124385	221895	2.59%	0.282%
50	19.169	2747	2751	2753	rBV	98593	147483	1.72%	0.188%
51	19.304	2770	2774	2788	rBV	295758	481323	5.63%	0.613%
52	19.522	2805	2811	2827	rBV	6301136	8492496	99.27%	10.808%
53	20.522	2978	2981	2985	rBV2	88889	115709	1.35%	0.147%
54	21.010	3060	3064	3072	rVB2	193041	351529	4.11%	0.447%
55	21.274	3105	3109	3118	rBV	4042238	5469481	63.93%	6.961%
56	23.504	3480	3488	3503	rVB	4175814	8555214	100.00%	10.888%
57	23.651	3507	3513	3530	rVB2	2589004	5357989	62.63%	6.819%

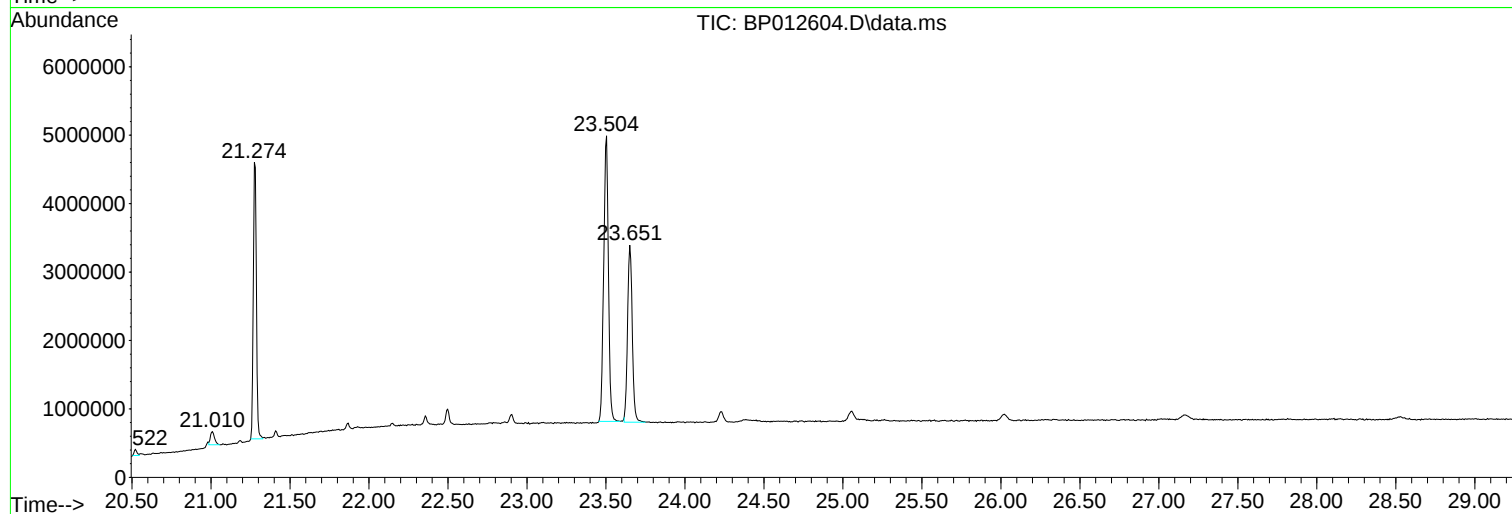
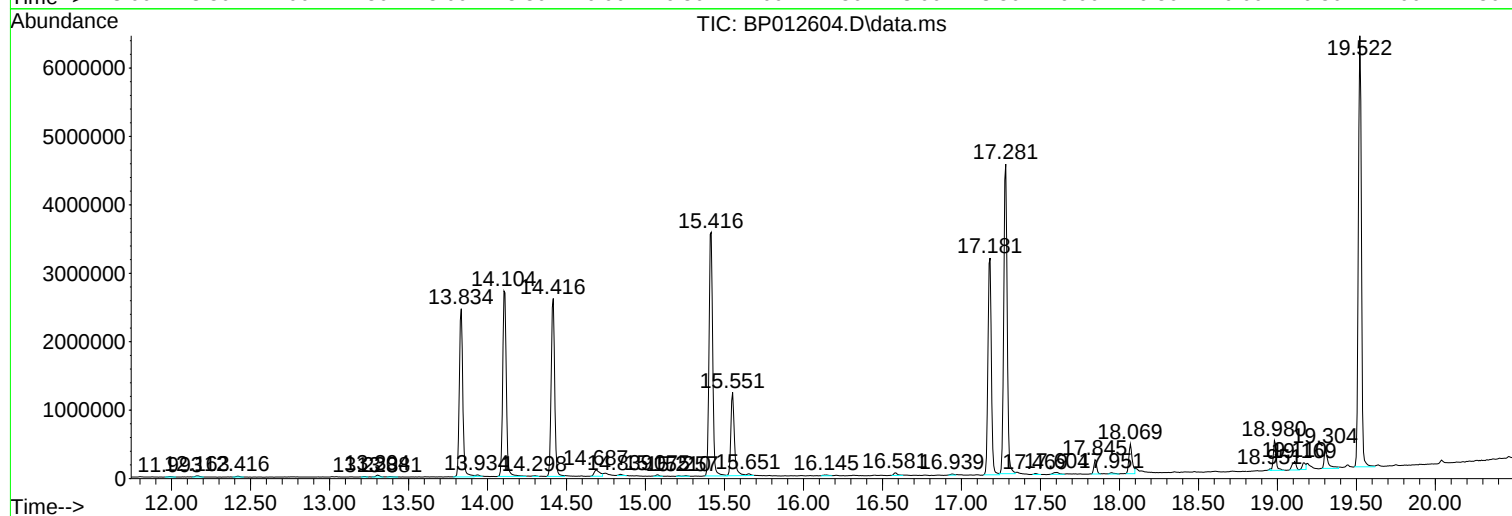
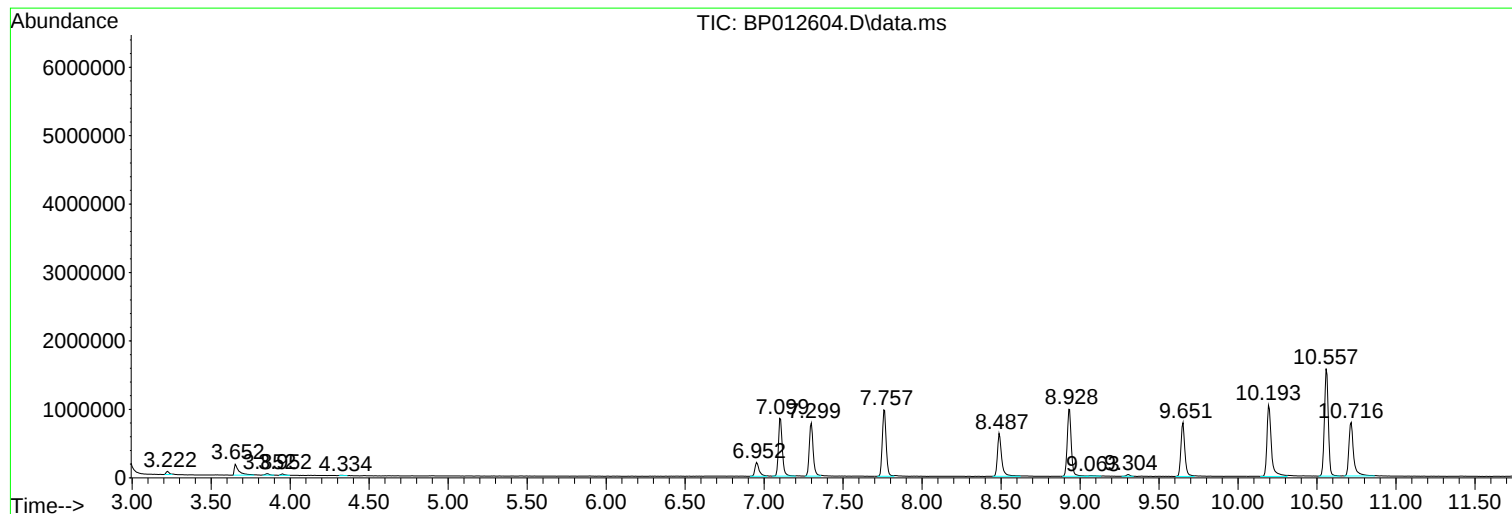
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110922.M
Quant Title : SVOA CALIBRATION

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



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Instrument :
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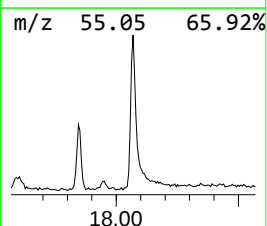
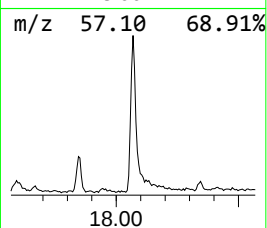
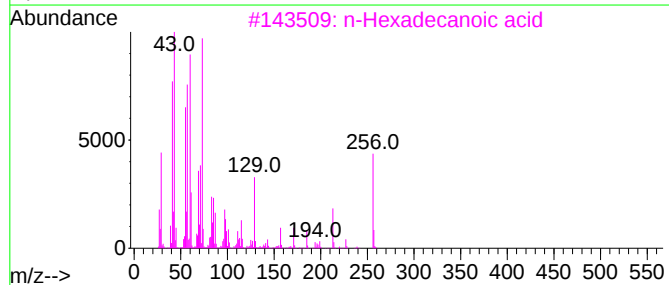
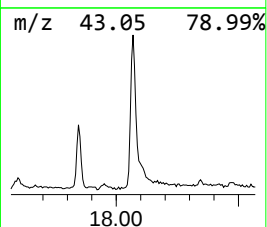
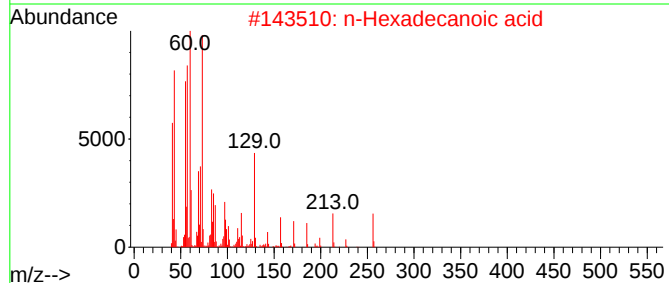
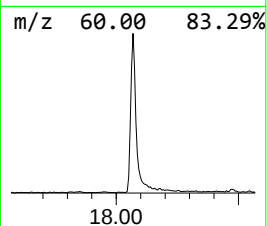
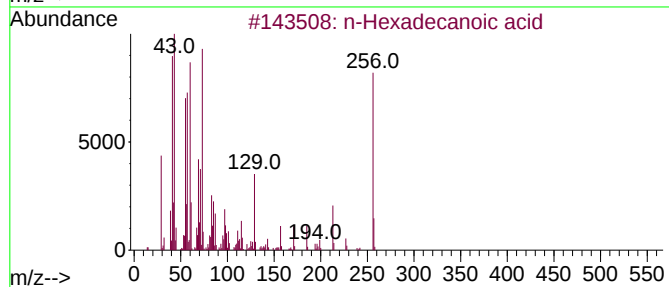
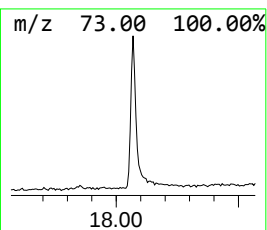
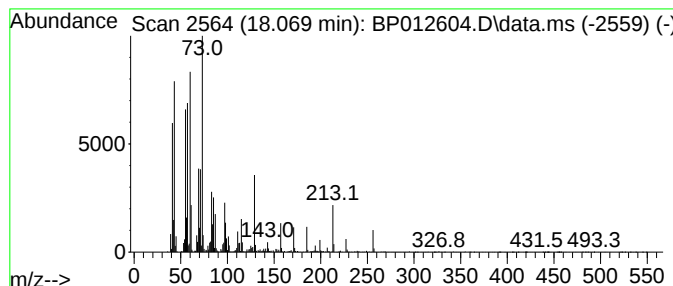
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110922.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 n-Hexadecanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.069	2.73 ng/ul	648206	Phenanthrene-d10	17.180

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	95
5			Tridecanoic acid	214	C13H26O2	000638-53-9	93



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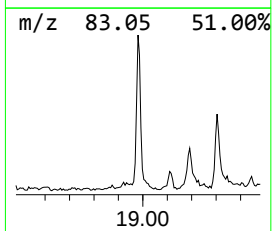
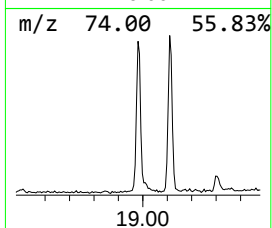
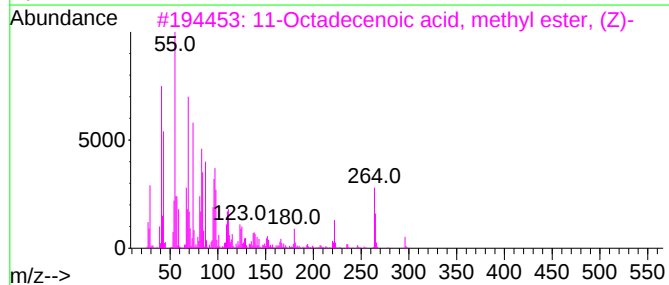
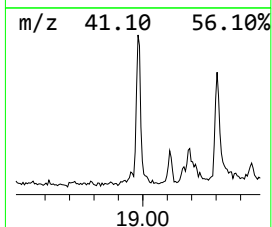
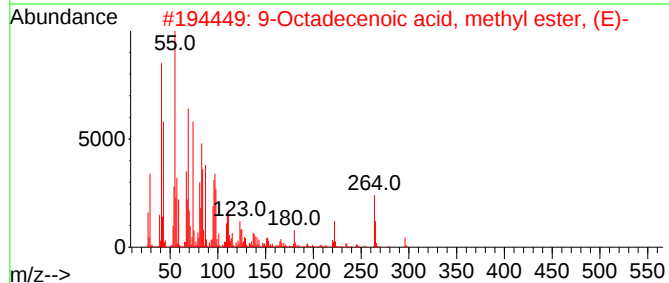
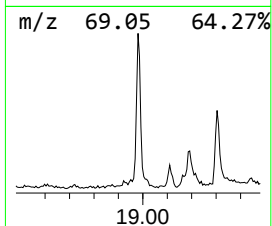
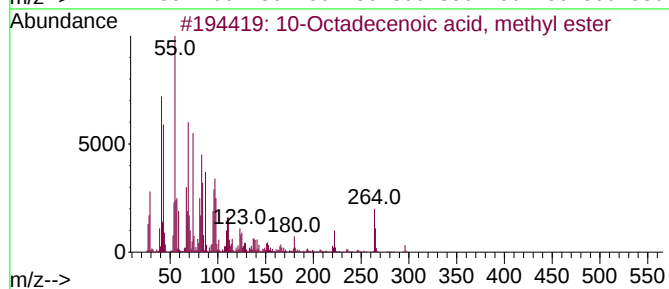
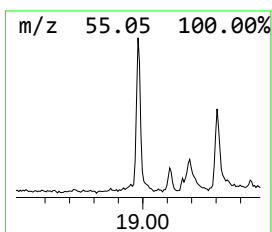
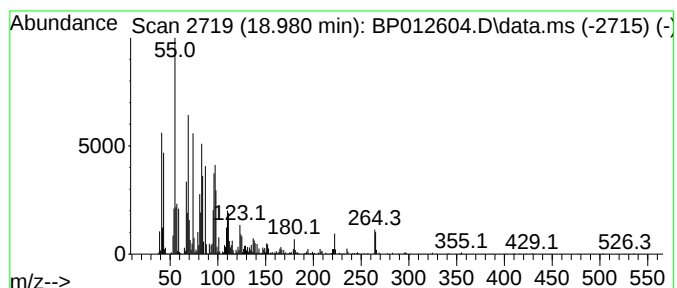
Instrument :
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Quant Title : SVOA CALIBRATION

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

Peak Number 2 10-Octadecenoic acid, methy... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.980	2.27 ng/ul	537510	Phenanthrene-d10	17.180	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	10-Octadecenoic acid, methyl ester	296	C19H36O2	013481-95-3	99
2	9-Octadecenoic acid, methyl ester	296	C19H36O2	001937-62-8	99
3	11-Octadecenoic acid, methyl ester	296	C19H36O2	001937-63-9	99
4	11-Octadecenoic acid, methyl ester	296	C19H36O2	052380-33-3	98
5	15-Octadecenoic acid, methyl ester	296	C19H36O2	004764-72-1	98



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TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
n-Hexadecanoic ...	18.069	2.7	ng/ul	648206	4	17.180	4742980	20.0
10-Octadecenoic...	18.980	2.3	ng/ul	537510	4	17.180	4742980	20.0