

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG112120\
 Data File : BG047408.D
 Acq On : 21 Nov 2020 23:52
 Operator : CG/JU
 Sample : L4711-17
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0AW9

Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : OFF
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 0.1 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG111920MA.M
 Title : SVOA CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.538	33	34	38	rBV2	2897	3297	0.25%	0.025%
2	3.614	40	47	52	rBV5	7580	14337	1.08%	0.110%
3	3.896	94	95	99	rVB2	2153	2348	0.18%	0.018%
4	4.049	117	121	124	rBV3	2602	3688	0.28%	0.028%
5	4.284	159	161	167	rVB3	3713	5116	0.39%	0.039%
6	4.354	170	173	175	rBV2	2745	2826	0.21%	0.022%
7	4.401	176	181	184	rVB2	2503	4140	0.31%	0.032%
8	4.437	184	187	188	rBV2	2158	2422	0.18%	0.019%
9	4.578	209	211	213	rBV	2282	2837	0.21%	0.022%
10	4.678	227	228	231	rBV	2534	2712	0.20%	0.021%
11	5.400	349	351	354	rBV	2128	3257	0.25%	0.025%
12	5.653	391	394	398	rVB2	1714	2381	0.18%	0.018%
13	6.581	550	552	556	rVB2	2612	2645	0.20%	0.020%
14	6.775	581	585	590	rBV2	2150	3590	0.27%	0.028%
15	7.410	682	693	703	rBV	185110	346140	26.15%	2.663%
16	7.504	706	709	712	rBV2				

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35	12.004	1471	1475	1476	rBV2	2249	2377	0.18%	0.018%
36	12.345	1529	1533	1537	rVB	2075	2253	0.17%	0.017%
37	12.662	1584	1587	1593	rVB	4778	7611	0.58%	0.059%
38	13.385	1706	1710	1714	rBV	2371	3238	0.24%	0.025%
39	13.497	1725	1729	1733	rBV2	3458	6013	0.45%	0.046%
40	13.726	1763	1768	1771	rBV3	4167	6363	0.48%	0.049%
41	13.767	1771	1775	1783	rVB3	27669	49414	3.73%	0.380%
42	13.820	1783	1784	1787	rBV3	2222	2366	0.18%	0.018%
43	14.278	1855	1862	1866	rBV	455428	678639	51.28%	5.220%
44	14.325	1866	1870	1876	rVB	163768	229564	17.35%	1.766%
45	14.507	1896	1901	1902	rBV	3171	3495	0.26%	0.027%
46	14.584	1907	1914	1921	rVV	468838	731761	55.29%	5.629%
47	14.666	1925	1928	1931	rVB	2859	2347	0.18%	0.018%
48	14.701	1931	1934	1936	rBV	2617	3172	0.24%	0.024%
49	14.889	1959	1966	1974	rVB	314551	516228	39.00%	3.971%
50	14.948	1974	1976	1979	rBV2				

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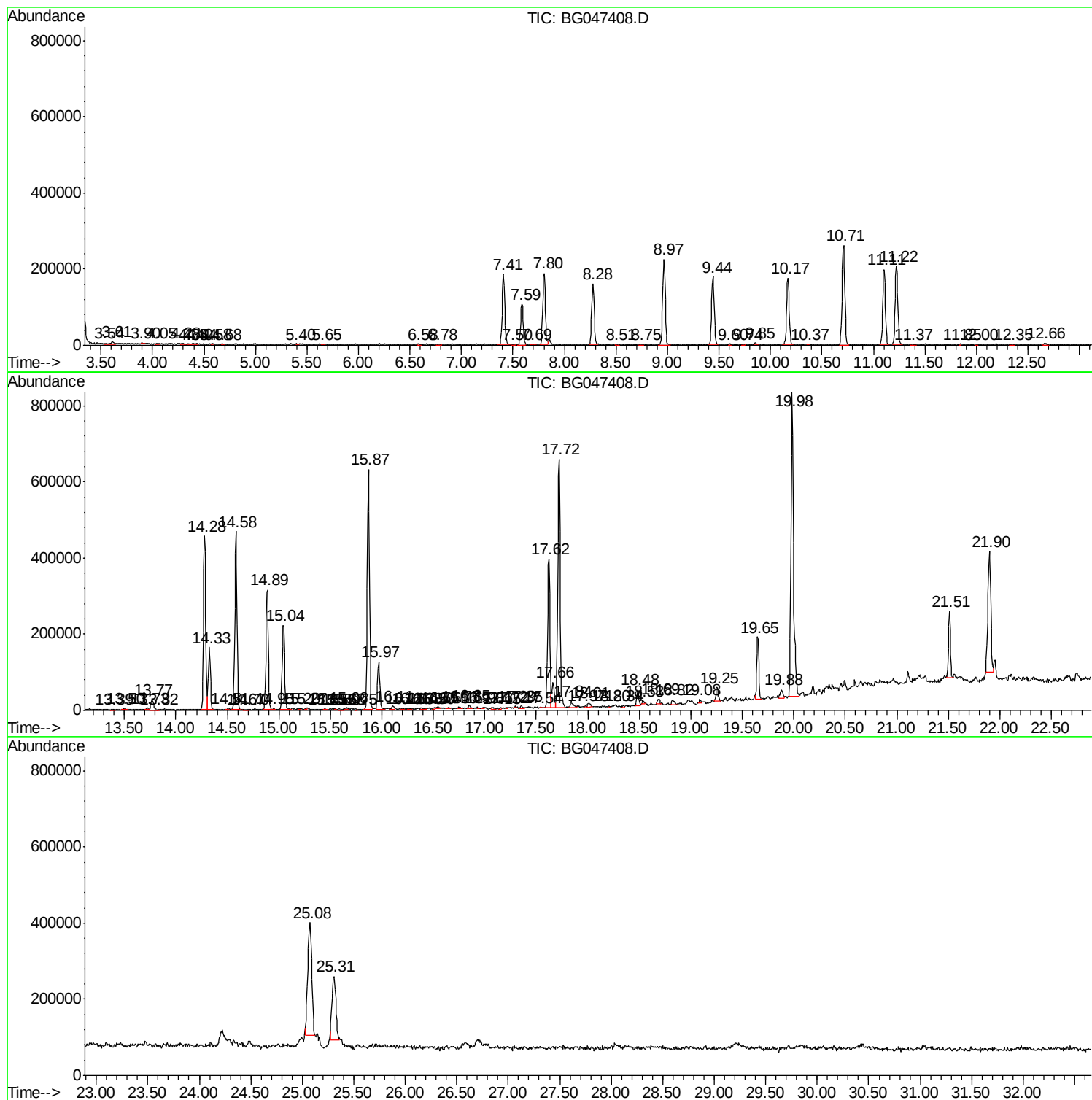
72	16.928	2312	2313	2315	rBV	4644	2718	0.21%	0.021%
73	16.981	2320	2322	2324	rBV3	3099	3000	0.23%	0.023%
74	17.063	2335	2336	2343	rVB4	4611	8293	0.63%	0.064%
75	17.151	2348	2351	2353	rBV3	2766	3636	0.27%	0.028%
76	17.216	2360	2362	2364	rBV3	3209	2370	0.18%	0.018%
77	17.292	2374	2375	2377	rBV2	4621	2608	0.20%	0.020%
78	17.351	2382	2385	2391	rBV6	5503	9616	0.73%	0.074%
79	17.545	2416	2418	2421	rBV4	2809	2371	0.18%	0.018%
80	17.621	2424	2431	2435	rVV	391123	617539	46.66%	4.750%
81	17.662	2435	2438	2442	rVV	67112	100859	7.62%	0.776%
82	17.721	2442	2448	2458	rVB	653236	1022867	77.29%	7.868%
83	17.839	2464	2468	2477	rBV5	17225	27859	2.10%	0.214%
84	17.974	2490	2491	2494	rVV2	4831	3789	0.29%	0.029%
85	18.009	2494	2497	2502	rVB4	11914	18903	1.43%	0.145%
86	18.203	2529	2530	2532	rBV	4790	2862	0.22%	0.022%
87	18.338	2552	2553	2556	rBV2	5749	6390	0.48%	0.049%

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG11920MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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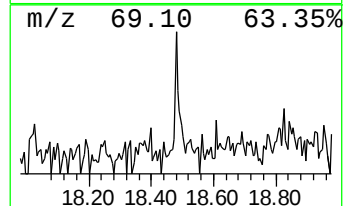
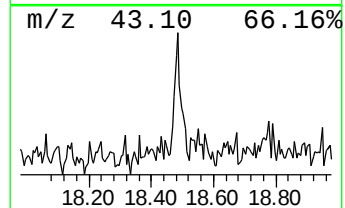
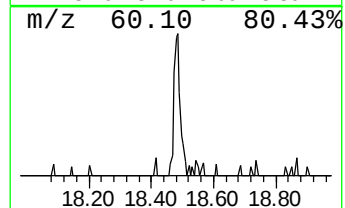
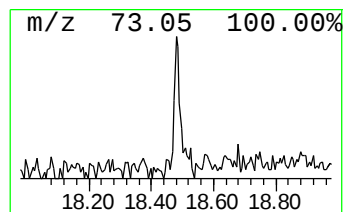
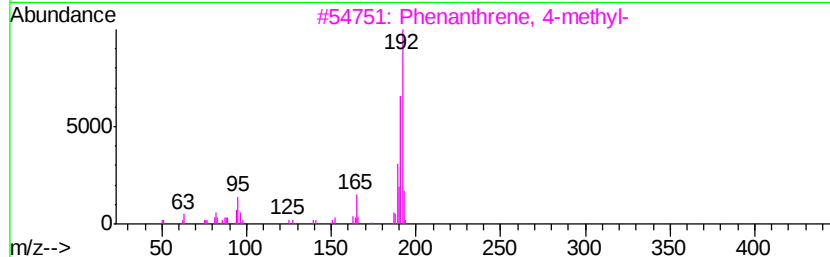
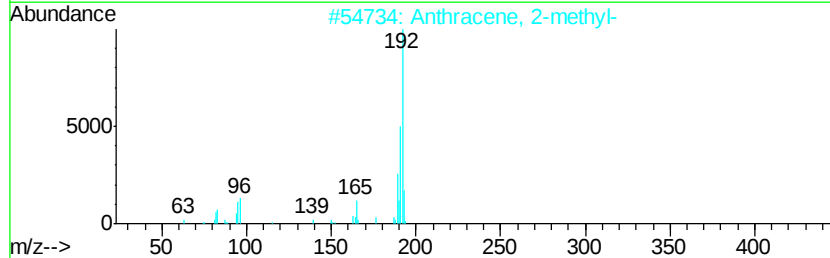
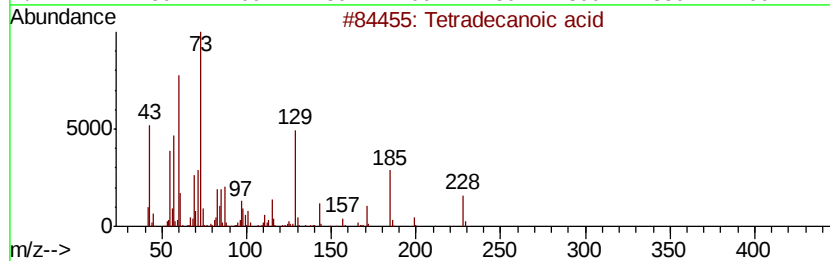
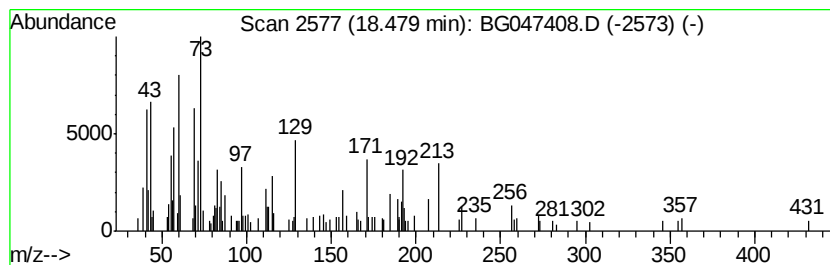
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Tetradecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.		
18.48	2.07 ng/ul	64032	Phenanthrene-d10	17.62		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tetradecanoic acid	228	C14H28O2	000544-63-8	38
2		Anthracene, 2-methyl-	192	C15H12	000613-12-7	25
3		Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	25
4		2-Hexyldecanoic acid	256	C16H32O2	025354-97-6	25
5		Anthracene, 9-methyl-	192	C15H12	000779-02-2	25



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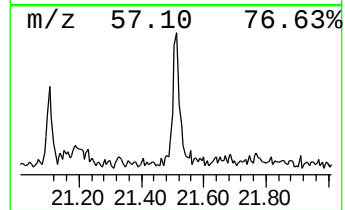
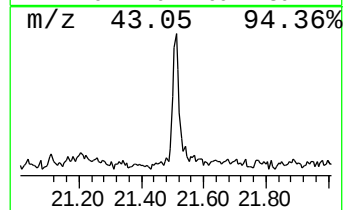
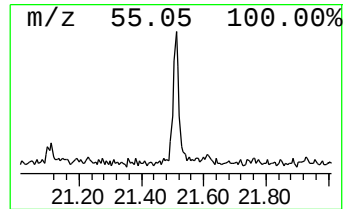
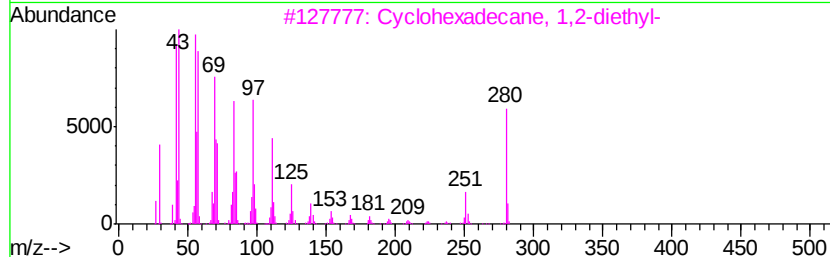
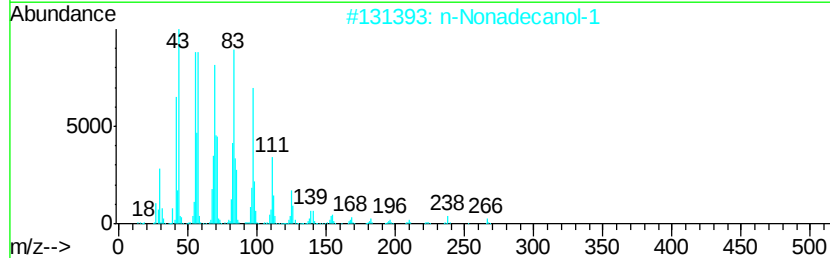
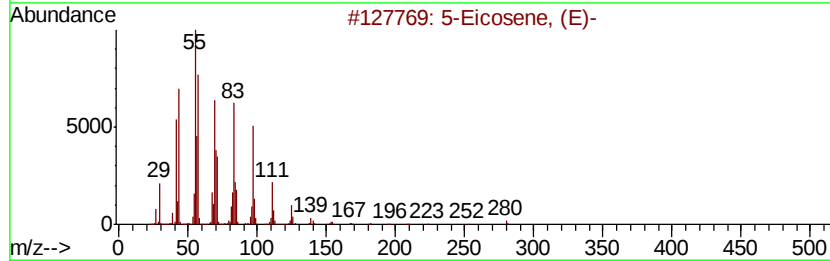
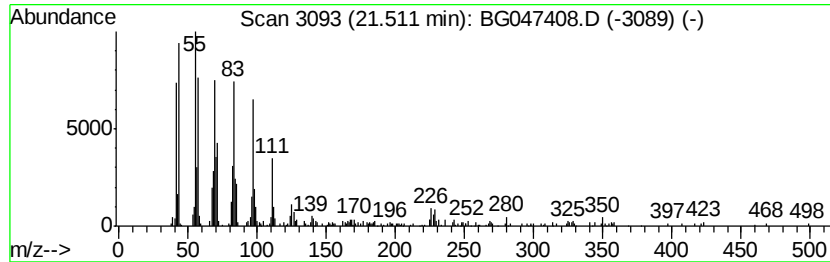
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.51	7.20 ng/ul	219441	Chrysene-d12	21.90

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Eicosene, (E)-	280	C20H40	074685-30-6	96
2		n-Nonadecanol-1	284	C19H40O	001454-84-8	95
3		Cyclohexadecane, 1,2-diethyl-	280	C20H40	1000155-85-3	94
4		Heptafluorobutyric acid, n-tetra...	410	C18H29F7O2	007365-36-8	94
5		Cycloeicosane	280	C20H40	000296-56-0	93



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TIC Top Hit name	RT	EstConc	Units	Response	---Internal Standard---			
					#	RT	Resp	Conc
Tetradecanoic acid	18.48	2.1	ng/ul	64032	4	17.62	617539	20.0
(DEL) Alkane: Str...	21.51	7.2	ng/ul	219441	5	21.90	609541	20.0