

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122823\
 Data File : BP019139.D
 Acq On : 28 Dec 2023 23:21
 Operator : MA/JU
 Sample : 06021-17
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BH7Z9

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-BP122723.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP019139.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.234	22	25	30	rVB	45136	51970	0.61%	0.092%
2	3.658	93	97	109	rBV	291465	429604	5.05%	0.759%
3	3.881	132	135	139	rVB2	9037	13064	0.15%	0.023%
4	3.975	148	151	157	rVB2	10778	14730	0.17%	0.026%
5	4.905	305	309	314	rBV	7589	11038	0.13%	0.020%
6	6.981	656	662	672	rVB	230104	366431	4.31%	0.648%
7	7.146	681	690	700	rBV	772909	1199772	14.11%	2.121%
8	7.334	715	722	732	rBV	753089	1190893	14.01%	2.105%
9	7.805	795	802	811	rBV	579107	903904	10.63%	1.598%
10	8.522	917	924	935	rBV	630235	1013297	11.92%	1.791%
11	8.969	992	1000	1012	rBV	883917	1419542	16.70%	2.509%
12	9.375	1064	1069	1075	rVB4	16094	29415	0.35%	0.052%
13	9.687	1113	1122	1132	rBV	685294	1127653	13.27%	1.993%
14	10.228	1205	1214	1229	rBV	1053070	1805965	21.25%	3.192%
15	10.599	1271	1277	1291	rVB	738723	1225851	14.42%	2.167%
16	10.746	1295	1302	1319	rBV	857970	1469352	17.29%	2.597%
17	11.869	1487	1493	1500	rVB2	17622	30119	0.35%	0.053%
18	12.198	1544	1549	1560	rVB	17820	35386	0.42%	0.063%
19	13.351	1738	1745	1751	rBV3	18570	31336	0.37%	0.055%
20	13.422	1753	1757	1763	rVB9	5417	9185	0.11%	0.016%
21	13.869	1825	1833	1843	rBV	2135582	2926845	34.43%	5.174%
22	14.140	1867	1879	1892	rBV	2382957	3433977	40.40%	6.070%
23	14.451	1922	1932	1939	rBV	1113166	1674363	19.70%	2.960%
24	14.669	1962	1969	1984	rBV	171415	333773	3.93%	0.590%
25	14.881	1999	2005	2009	rBV8	7175	13951	0.16%	0.025%
26	15.445	2094	2101	2109	rBV	3265000	4520908	53.18%	7.991%
27	15.569	2116	2122	2135	rBV	1009805	1384560	16.29%	2.447%
28	15.686	2138	2142	2149	rVB2	15238	24454	0.29%	0.043%
29	15.816	2158	2164	2169	rBV	126174	170069	2.00%	0.301%
30	16.228	2231	2234	2240	rVB4	13485	20658	0.24%	0.037%
31	16.381	2255	2260	2265	rBV	21029	28426	0.33%	0.050%
32	17.204	2394	2400	2410	rBV	1540740	2160564	25.42%	3.819%
33	17.304	2410	2417	2429	rVV2	3663729	5230373	61.53%	9.245%
34	17.392	2429	2432	2438	rVB6	12345	20810	0.24%	0.037%
35	18.098	2547	2552	2559	rBV	142719	191998	2.26%	0.339%
36	19.127	2722	2727	2732	rBV	55155	73816	0.87%	0.130%

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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

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37	19.198	2734	2739	2747	rBV	55277	94000	1.11%	0.166%
38	19.339	2759	2763	2769	rBV2	68314	93116	1.10%	0.165%
39	19.551	2793	2799	2806	rBV	5403114	6964336	81.93%	12.310%
40	21.027	3046	3050	3059	rBV3	182184	337089	3.97%	0.596%
41	21.310	3093	3098	3104	rBV	2181208	2856250	33.60%	5.049%
42	23.521	3467	3474	3488	rVB2	4427717	8500606	100.00%	15.026%
43	23.674	3493	3500	3511	rVB	1616883	3140145	36.94%	5.551%

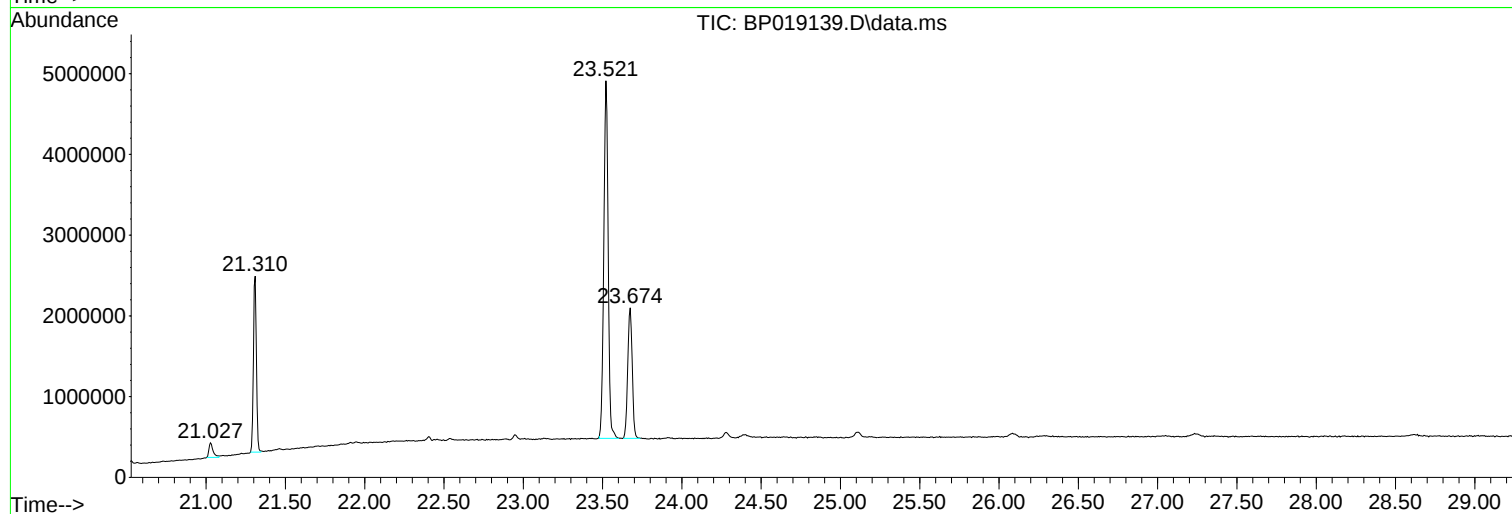
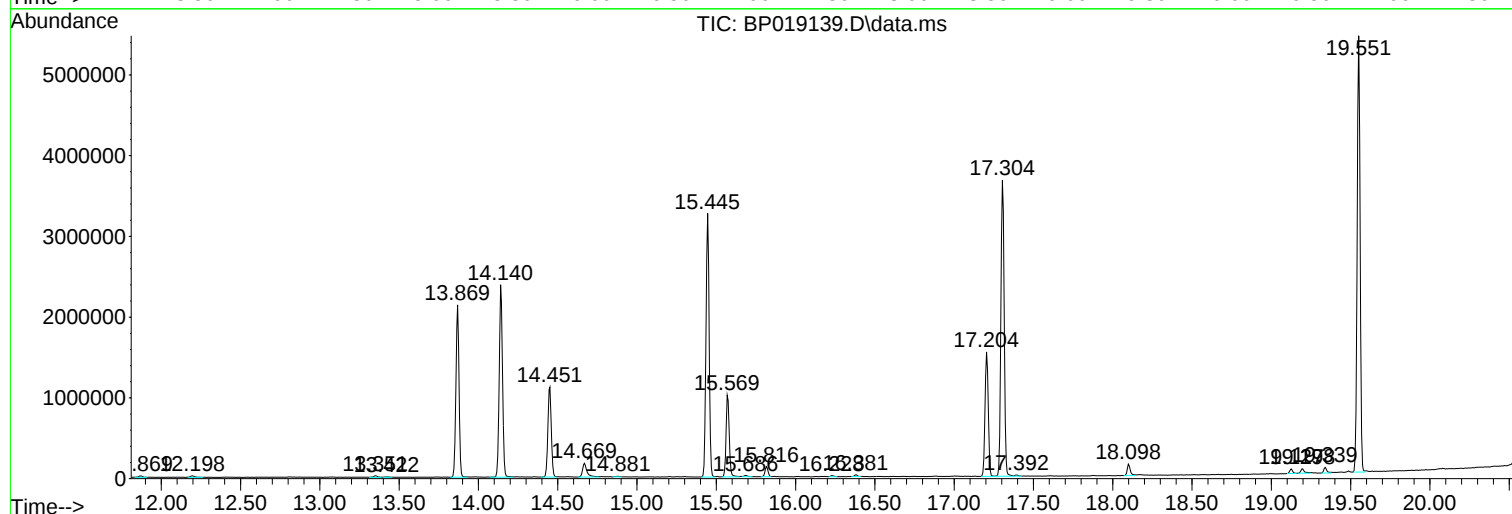
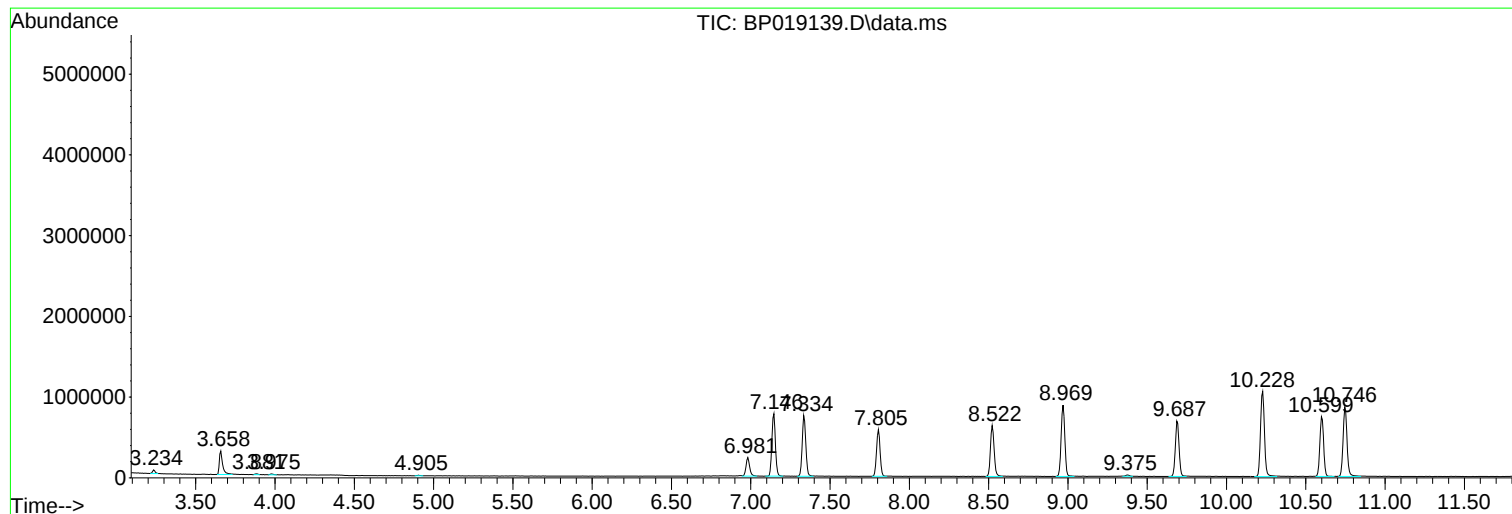
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP122723.MA.M
Quant Title : SVOA CALIBRATION

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



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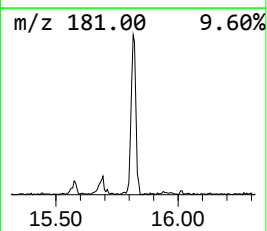
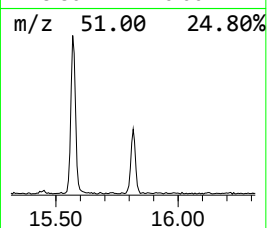
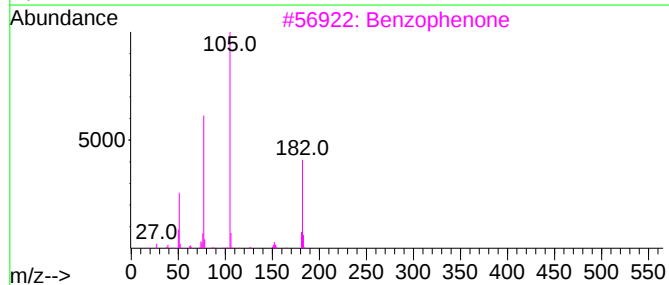
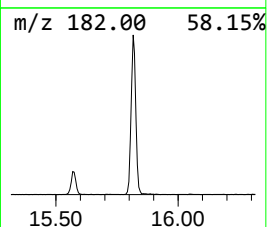
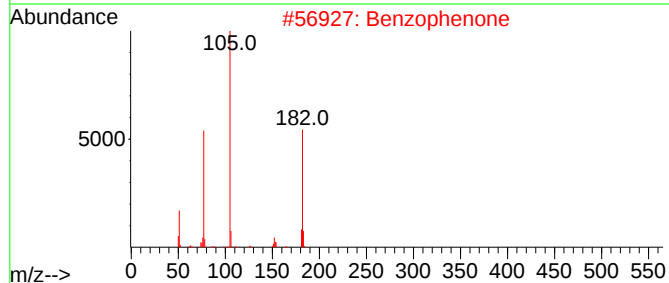
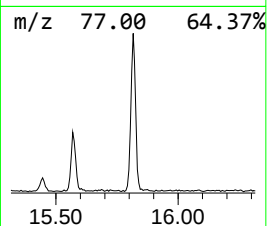
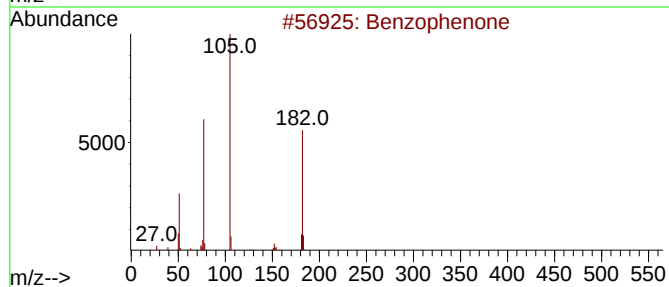
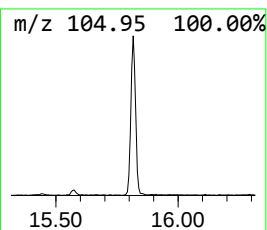
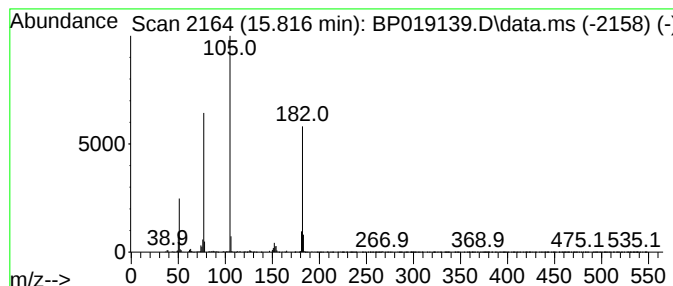
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP122723.MA.M
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TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Benzophenone Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.816	2.03 ng/ul	170069	Acenaphthene-d10	14.451

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	91
4			Benzophenone	182	C13H10O	000119-61-9	90
5			Azobenzene	182	C12H10N2	000103-33-3	64



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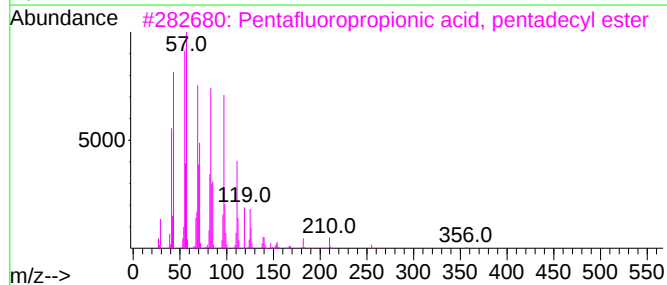
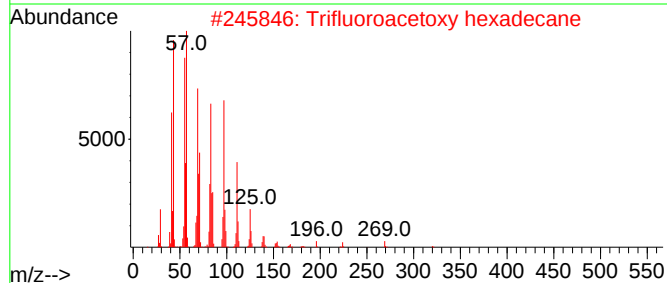
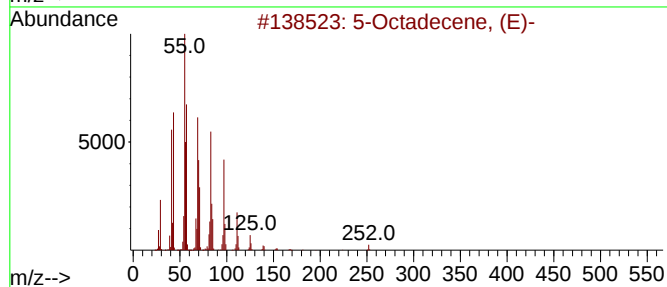
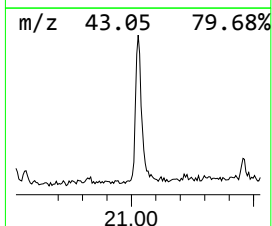
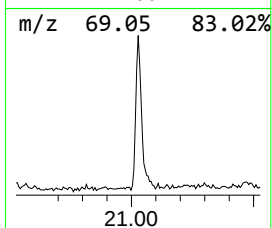
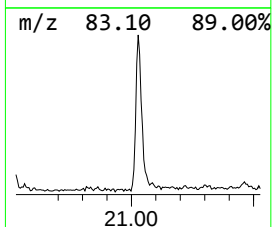
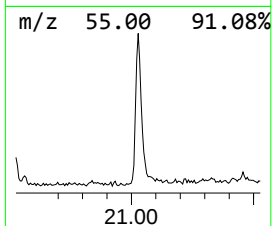
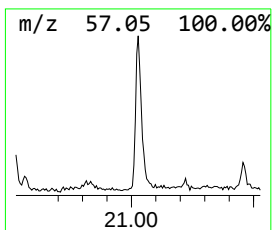
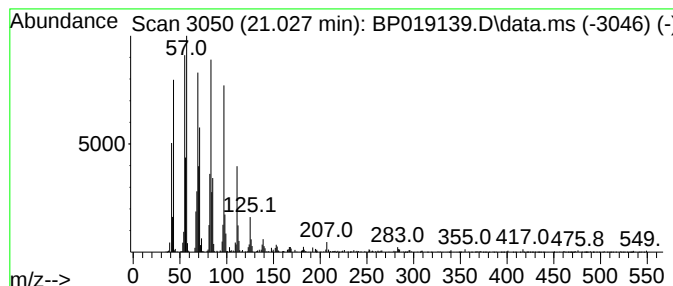
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Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.027	2.36 ng/ul	337089	Chrysene-d12	21.310

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			5-Octadecene, (E)-	252	C18H36	007206-21-5	94
2			Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	94
3			Pentafluoropropionic acid, penta...	374	C18H31F5O2	959092-08-1	94
4			Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	93
5			Acetic acid, chloro-, hexadecyl ...	318	C18H35ClO2	052132-58-8	91



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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzophenone	15.816	2.0	ng/u1	170069	3	14.451	1674360	20.0
(DEL) Alkane: S...	21.027	2.4	ng/u1	337089	5	21.310	2856250	20.0