

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121022\
 Data File : BP013044.D
 Acq On : 10 Dec 2022 17:35
 Operator : CG/JU
 Sample : N5950-02
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EYF49

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: BP013044.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.352	25	28	34	rBV	142613	182956	0.83%	0.098%
2	3.805	103	105	113	rBV	54128	98534	0.45%	0.053%
3	4.017	136	141	147	rVB	70840	106382	0.48%	0.057%
4	4.111	153	157	164	rVB	46123	70300	0.32%	0.038%
5	4.652	245	249	253	rBV5	16078	27932	0.13%	0.015%
6	7.128	664	670	679	rVB	647695	1046313	4.75%	0.560%
7	7.299	692	699	708	rBV	2452876	3842824	17.45%	2.055%
8	7.499	725	733	742	rBV	2236061	3477661	15.79%	1.860%
9	7.969	806	813	820	rBV	2539790	3973301	18.04%	2.125%
10	8.034	820	824	830	rVB	23637	36476	0.17%	0.020%
11	8.675	926	933	945	rBV	1839428	2926615	13.29%	1.565%
12	9.140	1004	1012	1026	rBV	3104748	5021875	22.80%	2.686%
13	9.299	1034	1039	1049	rVB	11000	24763	0.11%	0.013%
14	9.546	1075	1081	1093	rVB	176946	278834	1.27%	0.149%
15	9.863	1127	1135	1151	rBV	2397930	3935750	17.87%	2.105%
16	10.399	1218	1226	1240	rBV	3409683	5790287	26.29%	3.097%
17	10.705	1273	1278	1284	rBV3	26552	48757	0.22%	0.026%
18	10.787	1284	1292	1304	rVV	3691163	6165391	27.99%	3.297%
19	10.922	1308	1315	1328	rVV	673585	1188642	5.40%	0.636%
20	11.140	1348	1352	1359	rVB8	11564	22909	0.10%	0.012%
21	11.322	1377	1383	1388	rBV	25407	41845	0.19%	0.022%
22	11.381	1388	1393	1400	rVB2	15319	26928	0.12%	0.014%
23	11.799	1457	1464	1469	rBV9	12113	22165	0.10%	0.012%
24	12.040	1498	1505	1511	rBV2	66674	106638	0.48%	0.057%
25	12.193	1525	1531	1536	rVB2	55693	83750	0.38%	0.045%
26	12.363	1555	1560	1572	rVB	74377	127527	0.58%	0.068%
27	13.428	1736	1741	1747	rVB	83232	115115	0.52%	0.062%
28	13.499	1747	1753	1760	rVV	94644	156915	0.71%	0.084%
29	13.581	1761	1767	1773	rVB6	19610	43793	0.20%	0.023%
30	14.016	1834	1841	1851	rBV	7958438	11099509	50.40%	5.936%
31	14.134	1856	1861	1866	rVB9	24466	41348	0.19%	0.022%
32	14.304	1883	1890	1900	rBV	9000095	13133391	59.63%	7.024%
33	14.493	1917	1922	1927	rVB4	22439	35791	0.16%	0.019%
34	14.610	1935	1942	1952	rVB	6099158	8876753	40.30%	4.747%
35	14.798	1967	1974	1996	rBV	551709	953931	4.33%	0.510%
36	14.969	1996	2003	2006	rVV5	12542	22675	0.10%	0.012%

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

37	15.016	2006	2011	2017	rVB9	33243	64637	0.29%	0.035%
38	15.604	2103	2111	2124	rBV	11811352	17115051	77.71%	9.153%
39	15.722	2124	2131	2145	rVV	3924842	5437657	24.69%	2.908%
40	15.845	2147	2152	2158	rVB2	57761	97149	0.44%	0.052%
41	15.969	2167	2173	2179	rBV	1080173	1456322	6.61%	0.779%
42	16.387	2239	2244	2250	rVB8	17940	31858	0.14%	0.017%
43	16.745	2302	2305	2309	rBV5	22029	28008	0.13%	0.015%
44	17.157	2370	2375	2382	rVB4	20554	42890	0.19%	0.023%
45	17.363	2404	2410	2420	rBV	7525951	10604839	48.15%	5.672%
46	17.463	2420	2427	2444	rVV	13100749	19322294	87.73%	10.334%
47	17.645	2455	2458	2463	rVB	18668	29580	0.13%	0.016%
48	17.739	2470	2474	2481	rBV5	23394	31556	0.14%	0.017%
49	18.110	2534	2537	2542	rVV7	17475	26953	0.12%	0.014%
50	18.234	2553	2558	2567	rBV	404512	540302	2.45%	0.289%
51	18.310	2567	2571	2574	rVV	42902	53966	0.25%	0.029%
52	18.645	2625	2628	2631	rBV4	33950	36156	0.16%	0.019%
53	19.063	2696	2699	2702	rBV5	20186	28189	0.13%	0.015%
54	19.269	2730	2734	2739	rVB2	183235	240102	1.09%	0.128%
55	19.339	2741	2746	2749	rBV	177214	266691	1.21%	0.143%
56	19.463	2764	2767	2776	rVB5	88361	109307	0.50%	0.058%
57	19.610	2789	2792	2795	rVB2	68973	74176	0.34%	0.040%
58	19.698	2800	2807	2817	rBV	16160573	22024091	100.00%	11.779%
59	21.139	3048	3052	3063	rBV	876749	1374672	6.24%	0.735%
60	21.451	3099	3105	3115	rBV	7112319	9418784	42.77%	5.037%
61	22.722	3318	3321	3327	rVB	485145	671394	3.05%	0.359%
62	23.774	3493	3500	3513	rVB2	7821191	16587011	75.31%	8.871%
63	23.939	3521	3528	3540	rVB2	3904910	8111962	36.83%	4.338%

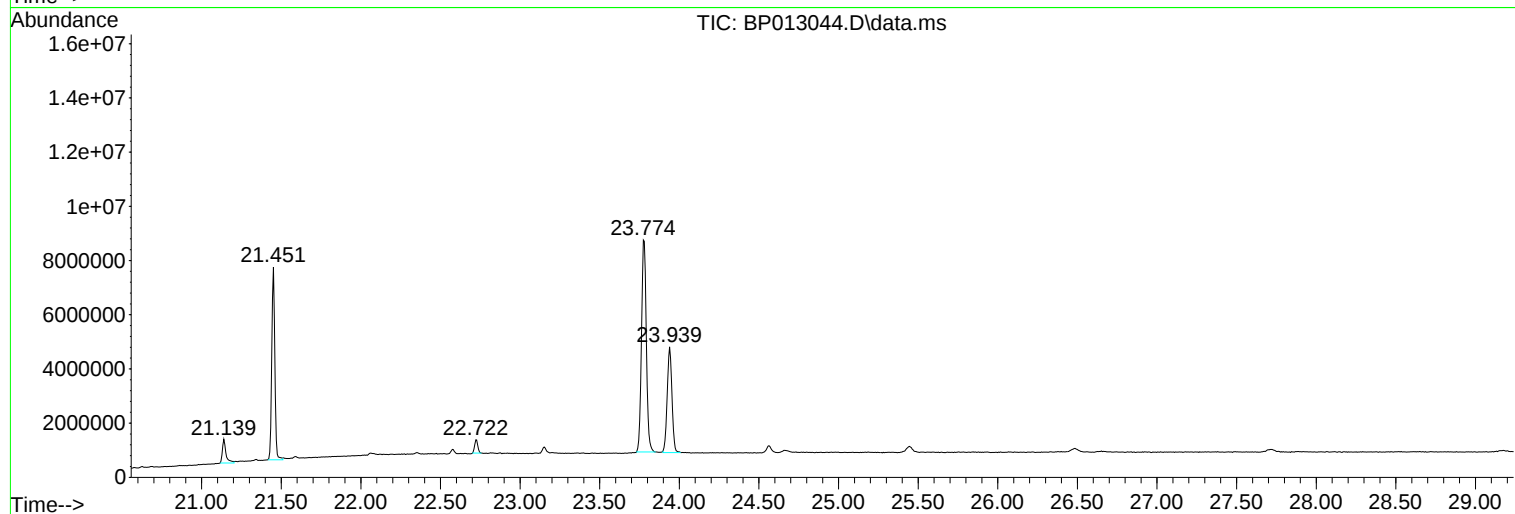
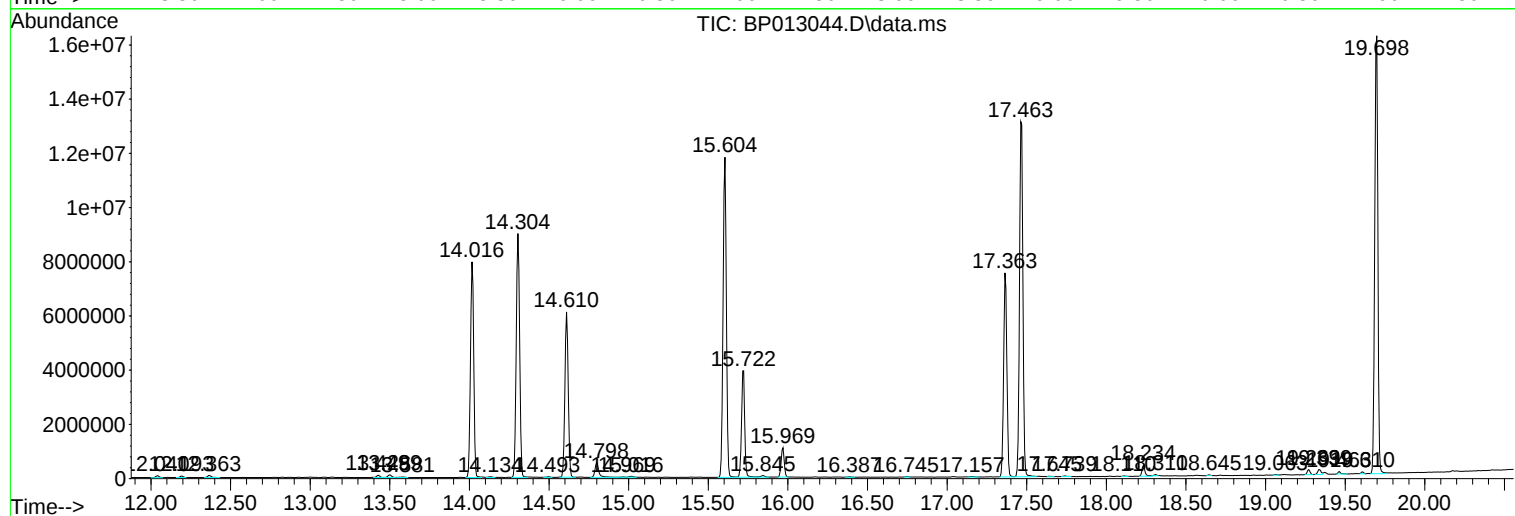
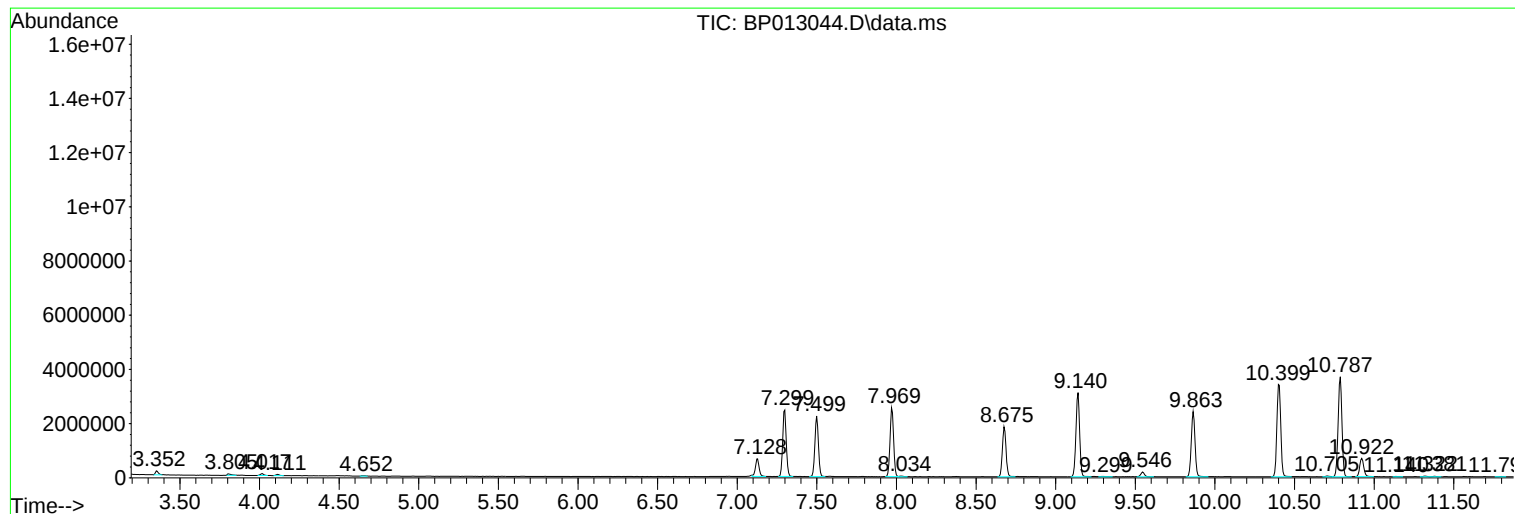
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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



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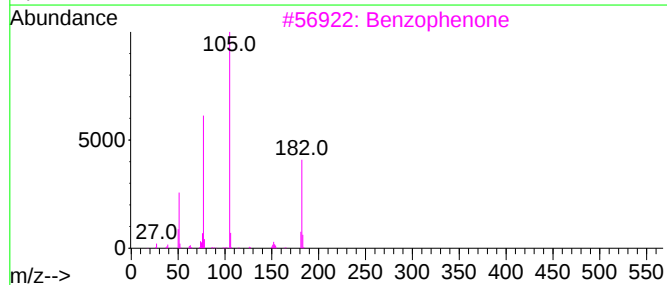
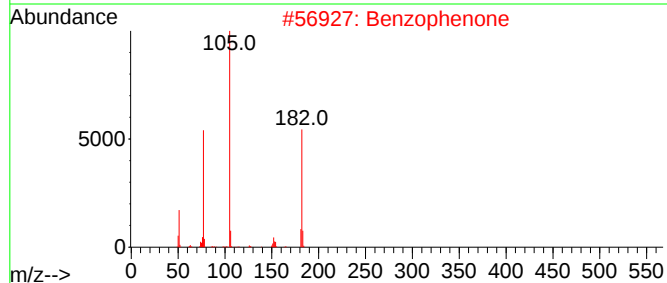
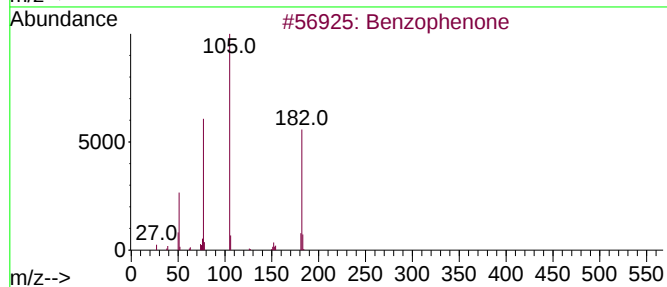
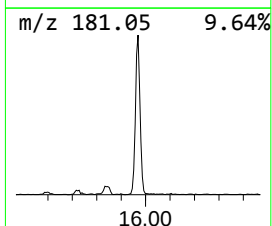
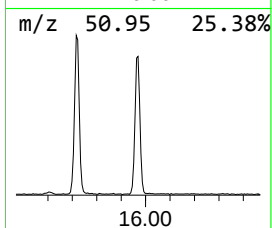
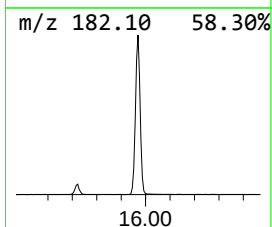
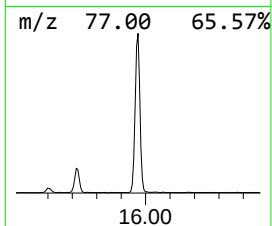
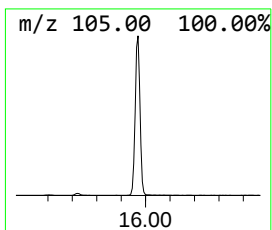
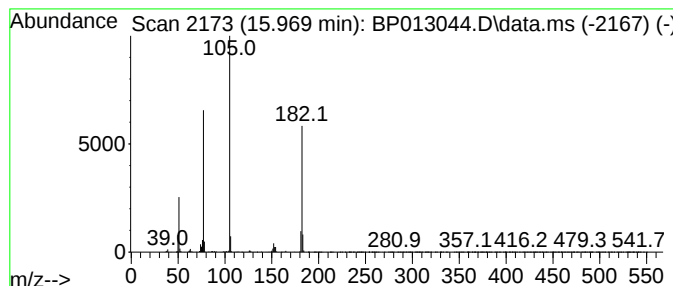
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzophenone Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.969	3.28 ng/ul	1456320	Acenaphthene-d10	14.610

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



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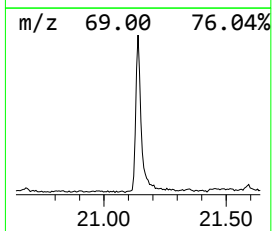
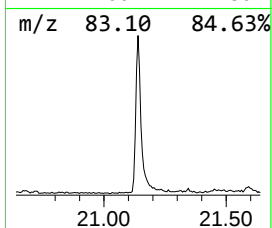
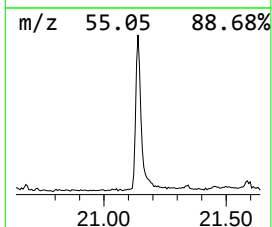
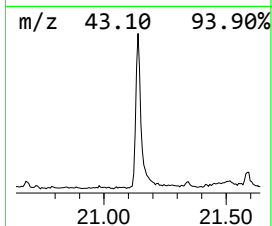
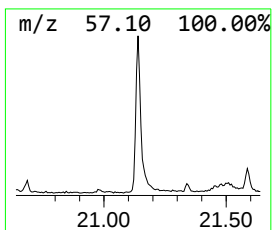
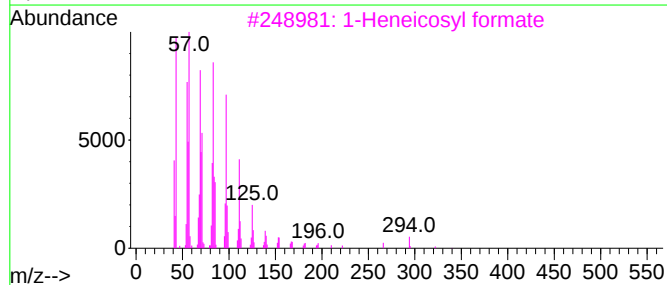
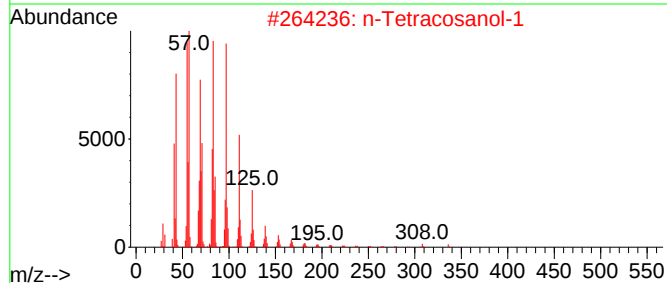
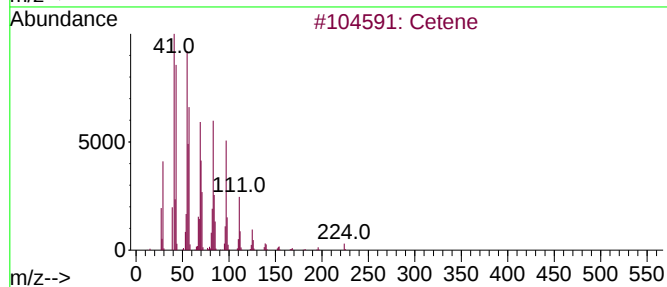
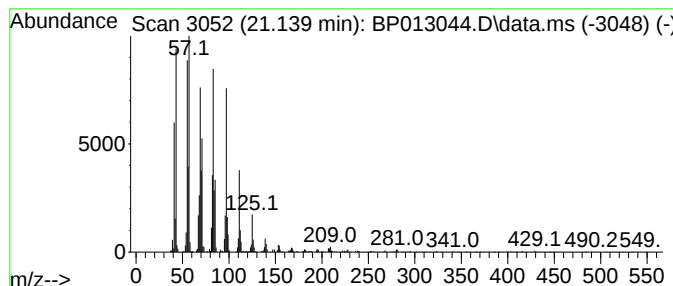
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.139	2.92 ng/ul	1374670	Chrysene-d12	21.451

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cetene	224	C16H32	000629-73-2	94
2			n-Tetracosanol-1	354	C24H50O	000506-51-4	94
3			1-Heneicosyl formate	340	C22H44O2	077899-03-7	94
4			Carbonic acid, octadecyl 2,2,2-t...	444	C21H39Cl3O3	1000314-56-3	94
5			Pentacos-1-ene	350	C25H50	016980-85-1	94



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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
Benzophenone	15.969	3.3	ng/u1	1456320	3	14.610	8876750	20.0
(DEL) Alkane: S...	21.139	2.9	ng/u1	1374670	5	21.451	9418780	20.0