

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112922\
 Data File : BP012855.D
 Acq On : 29 Nov 2022 16:55
 Operator : CG/JU
 Sample : N5683-01
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 ESQS8

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-BP112522.M
 Title : SVOA CALIBRATION

Signal : TIC: BP012855.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.393	31	35	42	rBV	94194	126079	0.75%	0.084%
2	3.823	104	108	122	rBV	666035	1076617	6.44%	0.721%
3	4.058	144	148	155	rVB	59095	90552	0.54%	0.061%
4	4.158	162	165	170	rVB	34478	41350	0.25%	0.028%
5	5.105	322	326	333	rVB5	10291	18165	0.11%	0.012%
6	7.158	668	675	688	rVB	466539	767265	4.59%	0.514%
7	7.334	696	705	714	rBV	1877722	2894624	17.31%	1.938%
8	7.540	732	740	748	rBV	1658278	2595339	15.52%	1.737%
9	7.628	751	755	760	rVB6	10159	18334	0.11%	0.012%
10	8.011	813	820	827	rBV	1690553	2681842	16.04%	1.795%
11	8.075	827	831	837	rVB4	14279	22344	0.13%	0.015%
12	8.710	932	939	948	rBV	1546802	2414915	14.44%	1.617%
13	9.081	997	1002	1006	rBV6	9384	18620	0.11%	0.012%
14	9.175	1010	1018	1030	rVV	1983533	3293494	19.70%	2.205%
15	9.275	1030	1035	1043	rVB2	32154	57642	0.34%	0.039%
16	9.599	1083	1090	1094	rBV2	41962	66035	0.39%	0.044%
17	9.905	1131	1142	1156	rBV	1575541	2559429	15.31%	1.713%
18	10.440	1224	1233	1250	rBV	2732191	4459947	26.68%	2.986%
19	10.599	1254	1260	1270	rBV	161443	285885	1.71%	0.191%
20	10.828	1291	1299	1306	rBV	2596492	4349088	26.01%	2.912%
21	10.957	1313	1321	1339	rBV	1886199	3334593	19.94%	2.232%
22	11.604	1424	1431	1434	rBV6	13975	24032	0.14%	0.016%
23	11.846	1465	1472	1476	rVB10	14032	20960	0.13%	0.014%
24	12.087	1507	1513	1518	rBV6	16302	32810	0.20%	0.022%
25	12.240	1533	1539	1543	rBV	34574	49461	0.30%	0.033%
26	12.410	1561	1568	1575	rBV	49952	82377	0.49%	0.055%
27	13.010	1665	1670	1673	rBV2	19590	30584	0.18%	0.020%
28	13.287	1714	1717	1721	rVB3	17921	24429	0.15%	0.016%
29	13.469	1743	1748	1755	rBV5	45207	85446	0.51%	0.057%
30	13.540	1755	1760	1764	rBV2	30462	44792	0.27%	0.030%
31	13.657	1777	1780	1787	rVB	28175	40123	0.24%	0.027%
32	14.051	1839	1847	1853	rBV	5817754	7809909	46.71%	5.228%
33	14.345	1889	1897	1905	rBV	6181202	9239267	55.26%	6.185%
34	14.540	1926	1930	1936	rVB8	19295	28852	0.17%	0.019%
35	14.645	1940	1948	1964	rBV	4182945	7137560	42.69%	4.778%
36	14.751	1964	1966	1973	rVV8	25644	47250	0.28%	0.032%

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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

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112522.M
 Title : SVOA CALIBRATION

37	14.828	1973	1979	1993	rVV	400980	663035	3.97%	0.444%
38	15.051	2013	2017	2023	rVB7	43040	71085	0.43%	0.048%
39	15.193	2037	2041	2046	rVB5	25665	38658	0.23%	0.026%
40	15.275	2050	2055	2062	rBV	106023	150552	0.90%	0.101%
41	15.381	2070	2073	2077	rBV5	16282	26550	0.16%	0.018%
42	15.469	2084	2088	2092	rBV3	50802	83592	0.50%	0.056%
43	15.640	2106	2117	2126	rBV	8418963	12348856	73.86%	8.267%
44	15.745	2126	2135	2152	rVV2	5796221	9962074	59.58%	6.669%
45	15.875	2155	2157	2165	rVB3	44796	76861	0.46%	0.051%
46	15.975	2169	2174	2175	rVV	138766	175620	1.05%	0.118%
47	16.004	2175	2179	2189	rVB	567315	825743	4.94%	0.553%
48	16.092	2189	2194	2210	rBV	318050	554901	3.32%	0.371%
49	16.245	2216	2220	2228	rVB	97878	142535	0.85%	0.095%
50	16.404	2242	2247	2253	rBV	250347	335065	2.00%	0.224%
51	16.787	2308	2312	2316	rBV6	24971	36424	0.22%	0.024%
52	17.404	2410	2417	2423	rBV	5305314	7610622	45.52%	5.095%
53	17.504	2427	2434	2445	rVV	9788192	14067665	84.14%	9.418%
54	17.586	2445	2448	2453	rVB3	47223	68580	0.41%	0.046%
55	18.269	2560	2564	2572	rBV	372178	515121	3.08%	0.345%
56	18.357	2574	2579	2587	rVB	379278	532310	3.18%	0.356%
57	19.304	2736	2740	2747	rBV2	132231	203242	1.22%	0.136%
58	19.369	2747	2751	2757	rBV	137795	262907	1.57%	0.176%
59	19.498	2769	2773	2781	rBV	250309	375137	2.24%	0.251%
60	19.645	2796	2798	2804	rVB5	52154	63191	0.38%	0.042%
61	19.728	2806	2812	2825	rVV	12326271	16719283	100.00%	11.193%
62	21.180	3055	3059	3064	rBV2	376980	517642	3.10%	0.347%
63	21.486	3106	3111	3122	rBV	5625015	7625408	45.61%	5.105%
64	23.839	3503	3511	3528	rBV	6213711	13007357	77.80%	8.708%
65	23.998	3531	3538	3551	rVB2	3064452	6446065	38.55%	4.315%

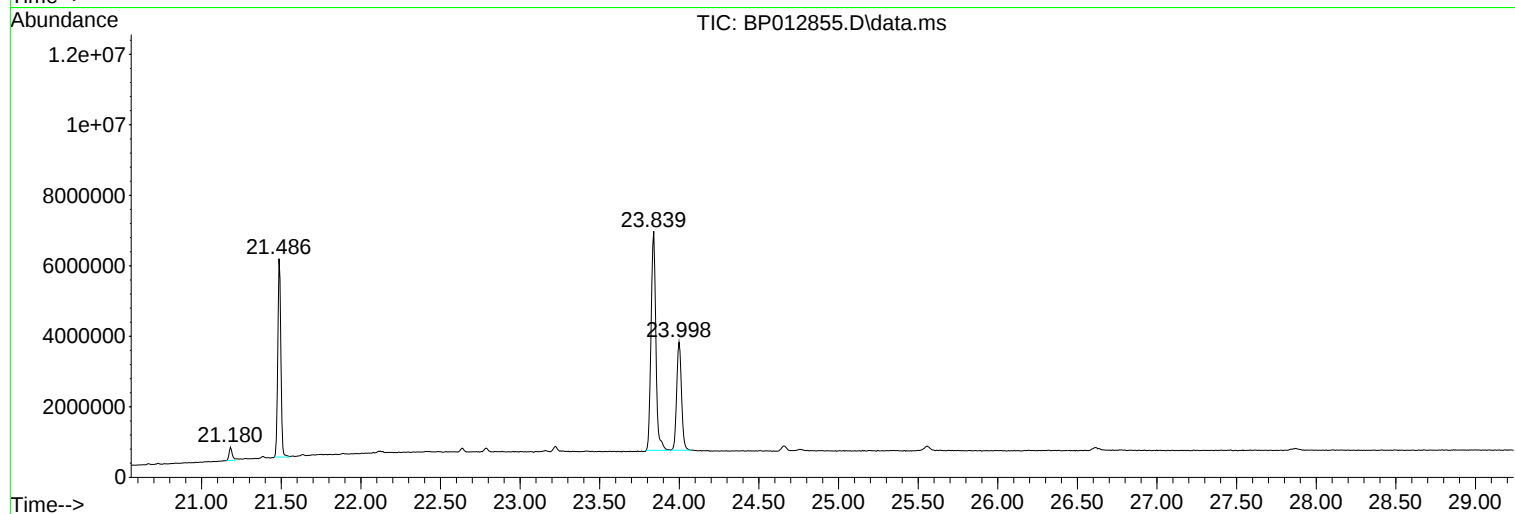
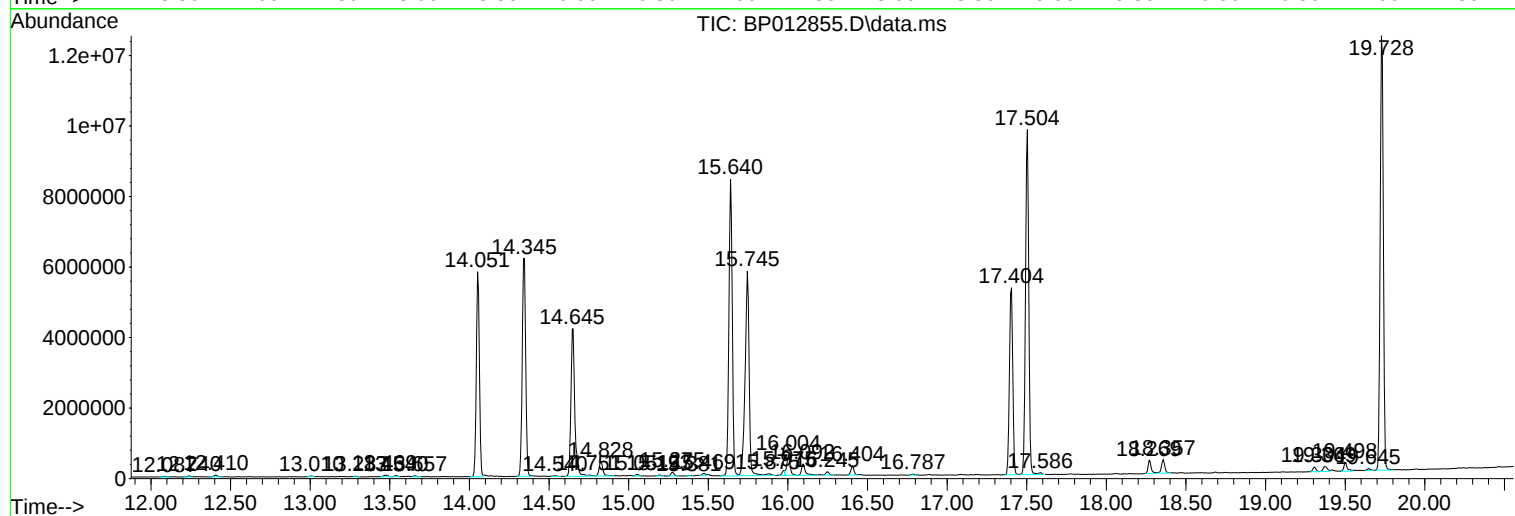
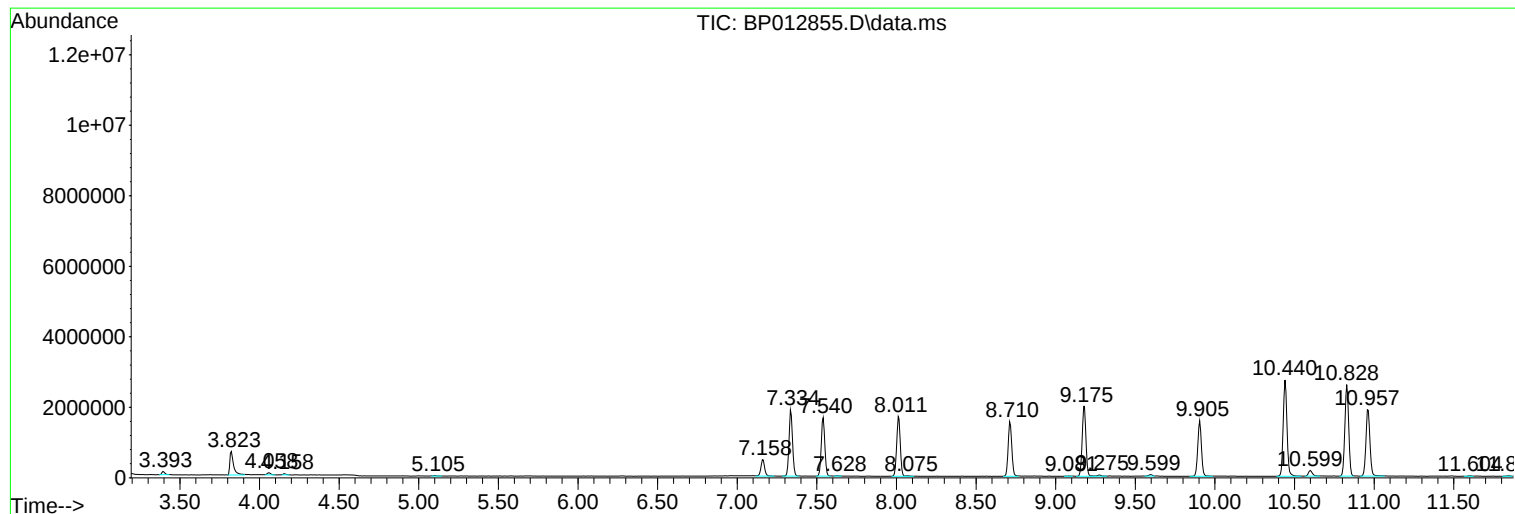
Sum of corrected areas: 149376092

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112522.M
Quant Title : SVOA CALIBRATION

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



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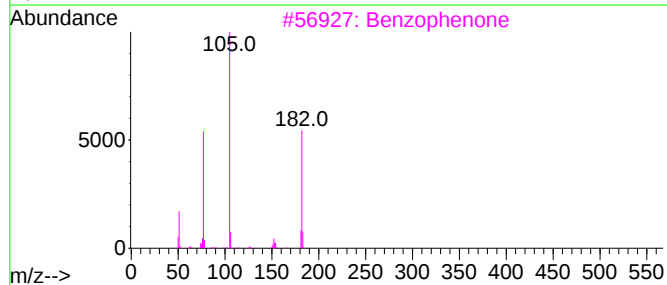
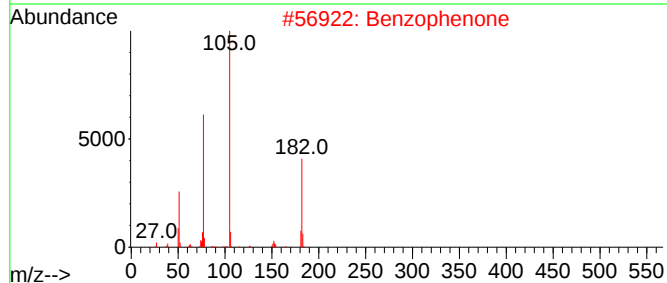
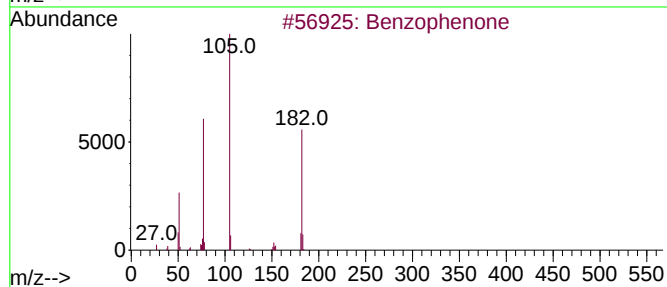
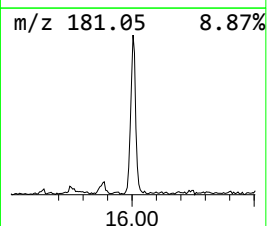
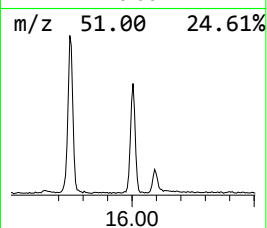
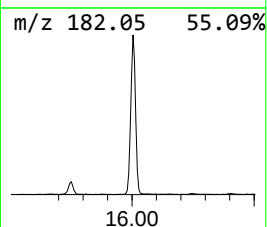
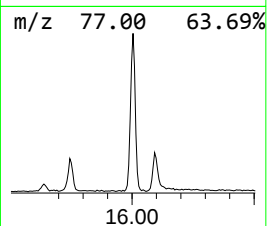
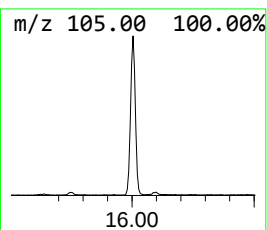
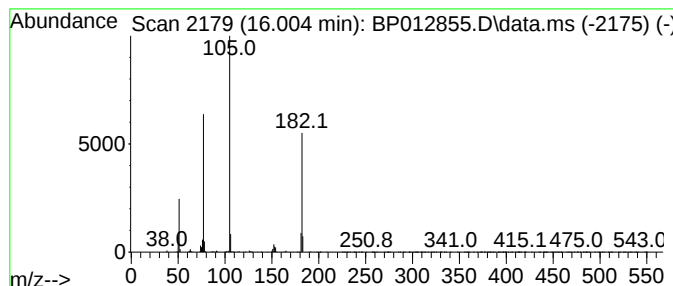
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112522.M
Quant Title : SVOA CALIBRATION

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.
16.004	2.31 ng/ul	825743	Acenaphthene-d10	14.646

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	90
5			1,2,4,5-Tetroxane, 3,3,6,6-tetra...	396	C26H20O4	016204-36-7	78



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzophenone	16.004	2.3	ng/u1	825743	3	14.646	7137560	20.0