

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051723\
 Data File : BP015126.D
 Acq On : 17 May 2023 14:58
 Operator : CG/JU
 Sample : 02768-16
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 CONP6

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-BP051123.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP015126.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.434	4	8	14	rVB	68946	98274	0.91%	0.104%
2	3.875	81	83	104	rBV	330248	576666	5.32%	0.608%
3	4.111	118	123	131	rVB	34742	61726	0.57%	0.065%
4	4.217	137	141	146	rVB	16027	24127	0.22%	0.025%
5	4.599	204	206	212	rVV7	9496	13166	0.12%	0.014%
6	4.675	215	219	226	rVB	48517	68495	0.63%	0.072%
7	5.169	299	303	312	rVB5	7926	15528	0.14%	0.016%
8	5.664	382	387	392	rBV	12164	19446	0.18%	0.021%
9	5.787	403	408	414	rVB	26208	40421	0.37%	0.043%
10	6.146	464	469	479	rBV	29680	48636	0.45%	0.051%
11	7.234	648	654	657	rBV2	35445	55817	0.51%	0.059%
12	7.275	657	661	677	rVB	290226	478951	4.42%	0.505%
13	7.446	683	690	702	rBV	1332755	2173843	20.04%	2.292%
14	7.652	717	725	736	rBV	1204820	1942776	17.91%	2.048%
15	7.852	755	759	766	rVB6	7027	13305	0.12%	0.014%
16	8.128	797	806	814	rBV	964707	1587881	14.64%	1.674%
17	8.187	814	816	823	rVB8	8185	14110	0.13%	0.015%
18	8.840	917	927	938	rBV	963734	1596565	14.72%	1.683%
19	8.952	941	946	952	rVB9	5533	14330	0.13%	0.015%
20	9.310	998	1007	1018	rBV	1580358	2667689	24.59%	2.813%
21	9.410	1018	1024	1031	rVV	50468	82797	0.76%	0.087%
22	9.704	1067	1074	1081	rVB	120296	180009	1.66%	0.190%
23	10.034	1122	1130	1146	rBV	1287591	2121766	19.56%	2.237%
24	10.575	1214	1222	1237	rBV	1876992	3131361	28.87%	3.301%
25	10.963	1280	1288	1302	rVB	1594300	2614379	24.10%	2.756%
26	11.099	1302	1311	1329	rBV	1670474	2915280	26.88%	3.074%
27	11.746	1415	1421	1424	rBV7	10222	19831	0.18%	0.021%
28	11.769	1424	1425	1431	rVB6	6567	11020	0.10%	0.012%
29	12.198	1493	1498	1504	rVB	78706	118365	1.09%	0.125%
30	12.546	1550	1557	1569	rBV3	32409	85680	0.79%	0.090%
31	13.040	1637	1641	1649	rBV8	11701	20566	0.19%	0.022%
32	13.657	1741	1746	1752	rVB4	12528	21207	0.20%	0.022%
33	14.181	1828	1835	1848	rVV	3926720	5442107	50.17%	5.738%
34	14.475	1873	1885	1895	rBV	4429333	6584577	60.71%	6.942%
35	14.557	1895	1899	1908	rVB	27314	51259	0.47%	0.054%
36	14.775	1929	1936	1945	rBV	2456343	3622669	33.40%	3.819%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051123.MA.M
 Title : SVOA CALIBRATION

37	14.963	1962	1968	1982	rBV	227424	456130	4.21%	0.481%
38	15.187	2001	2006	2012	rVB9	18054	33566	0.31%	0.035%
39	15.510	2056	2061	2065	rBV2	15253	26497	0.24%	0.028%
40	15.728	2090	2098	2099	rBV	65962	89522	0.83%	0.094%
41	15.769	2099	2105	2118	rVV	6025000	8584084	79.14%	9.050%
42	15.887	2118	2125	2140	rVV	1992920	2789230	25.71%	2.941%
43	16.004	2142	2145	2152	rVB2	27415	45065	0.42%	0.048%
44	16.128	2161	2166	2181	rVB	326847	473982	4.37%	0.500%
45	16.469	2215	2224	2227	rBV7	12911	27599	0.25%	0.029%
46	16.981	2306	2311	2315	rBV	60134	83812	0.77%	0.088%
47	17.204	2346	2349	2355	rVB8	8296	13156	0.12%	0.014%
48	17.322	2365	2369	2373	rVV5	10723	18766	0.17%	0.020%
49	17.533	2398	2405	2415	rBV	3082485	4331719	39.94%	4.567%
50	17.633	2415	2422	2442	rVV	6733029	9508287	87.66%	10.025%
51	17.804	2448	2451	2457	rVB3	12755	15739	0.15%	0.017%
52	18.045	2488	2492	2497	rBV	48422	61844	0.57%	0.065%
53	18.175	2511	2514	2519	rVB6	12085	16215	0.15%	0.017%
54	18.386	2545	2550	2558	rBV	116384	162111	1.49%	0.171%
55	18.463	2559	2563	2571	rVB	36406	53457	0.49%	0.056%
56	18.951	2643	2646	2649	rVB	43902	43385	0.40%	0.046%
57	19.428	2723	2727	2733	rBV2	88633	111537	1.03%	0.118%
58	19.498	2734	2739	2744	rVV	85588	123433	1.14%	0.130%
59	19.610	2755	2758	2764	rBV3	54462	78730	0.73%	0.083%
60	19.716	2773	2776	2779	rBV2	39358	43033	0.40%	0.045%
61	19.851	2793	2799	2808	rBV	8544481	10846771	100.00%	11.436%
62	20.386	2888	2890	2894	rVB2	43148	45916	0.42%	0.048%
63	21.286	3040	3043	3051	rBV4	120053	222954	2.06%	0.235%
64	21.610	3093	3098	3111	rVB	3288036	4409942	40.66%	4.649%
65	24.015	3499	3507	3527	rBV2	4274996	9637077	88.85%	10.160%
66	24.180	3529	3535	3552	rVB	1859869	3962180	36.53%	4.177%

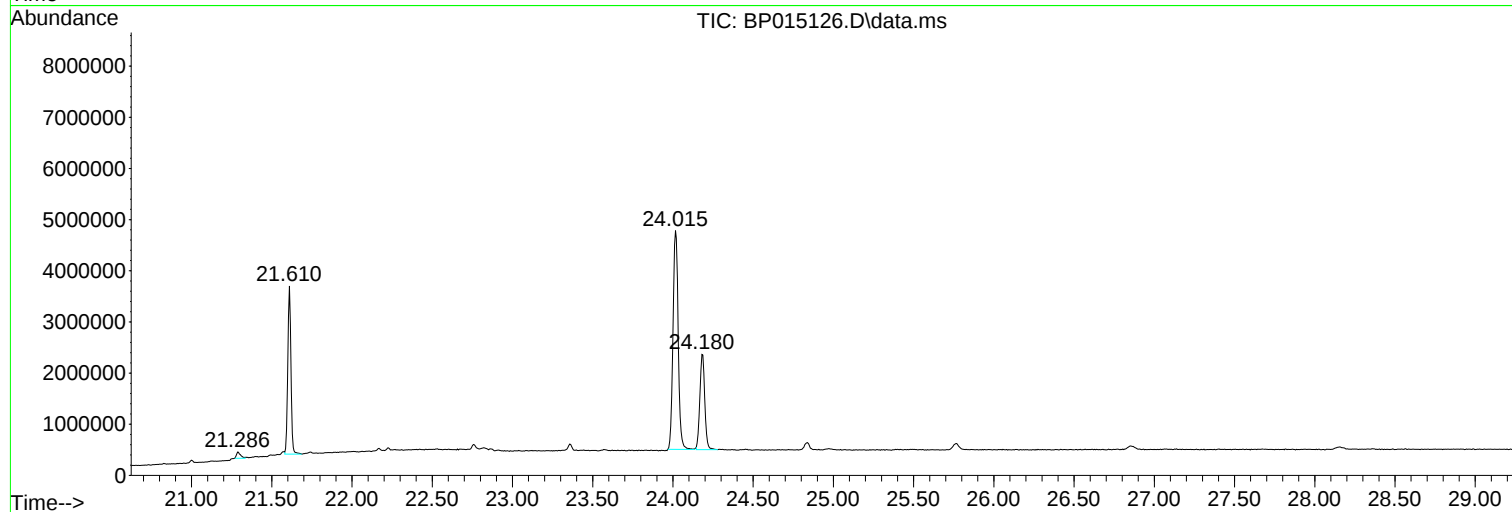
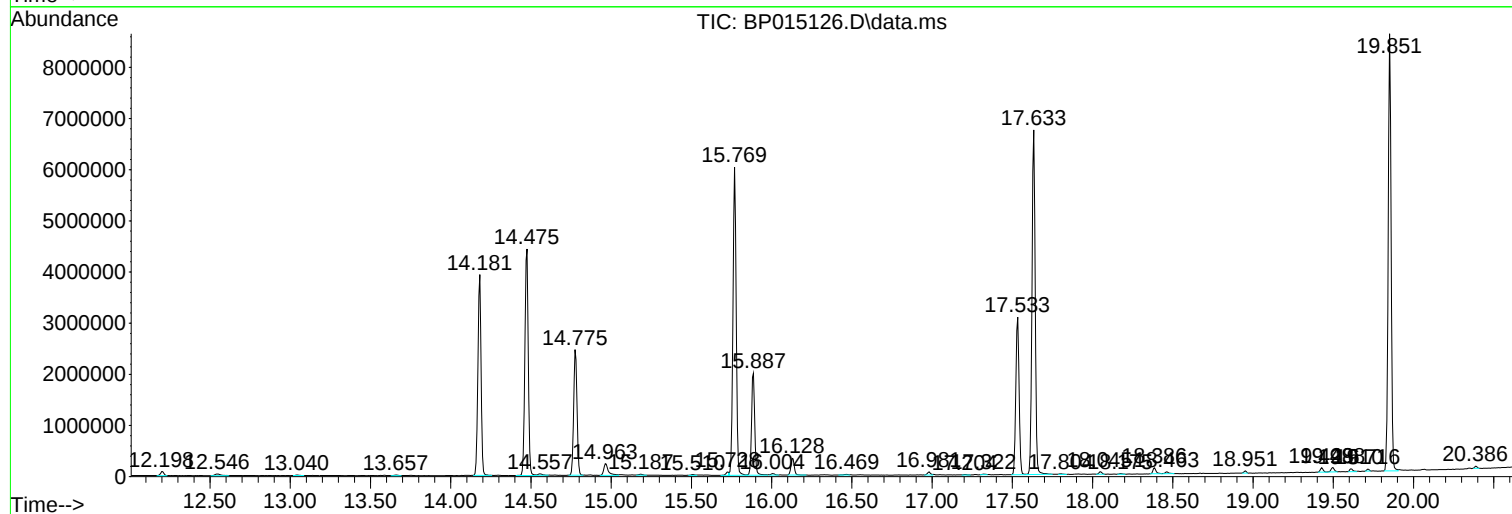
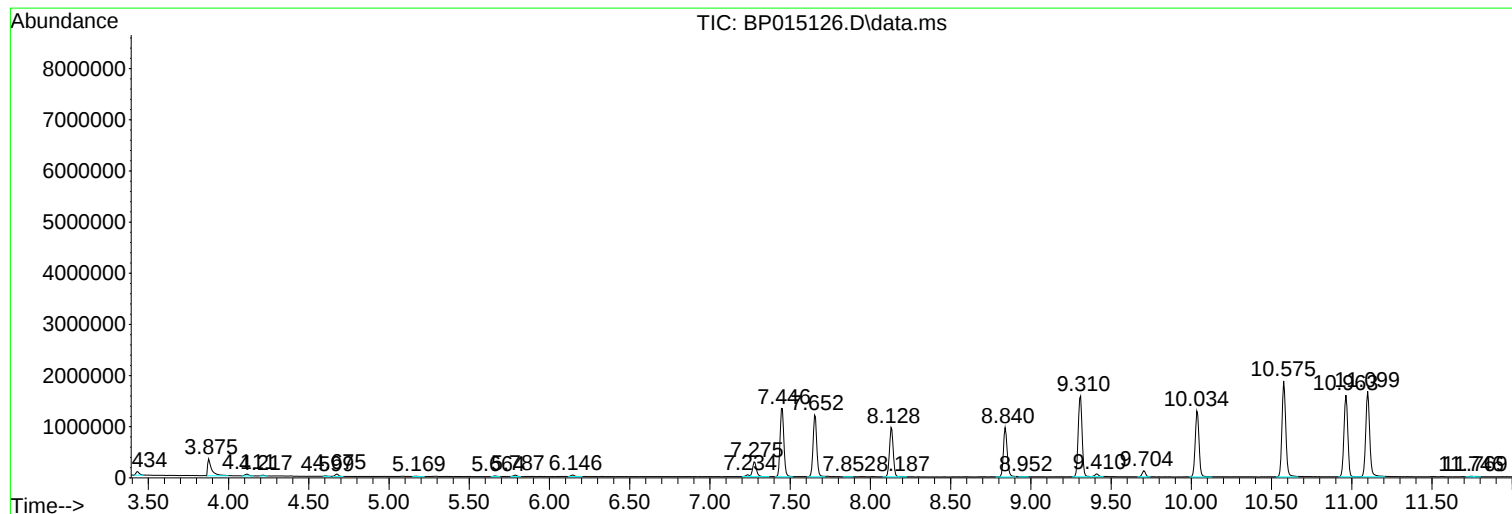
Sum of corrected areas: 94850364

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051123.MA.M
Quant Title : SVOA CALIBRATION

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



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Instrument :
BNA_P
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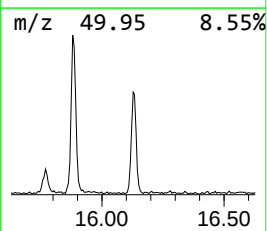
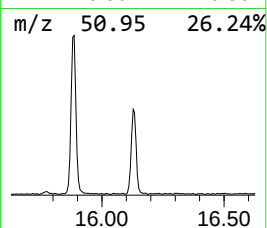
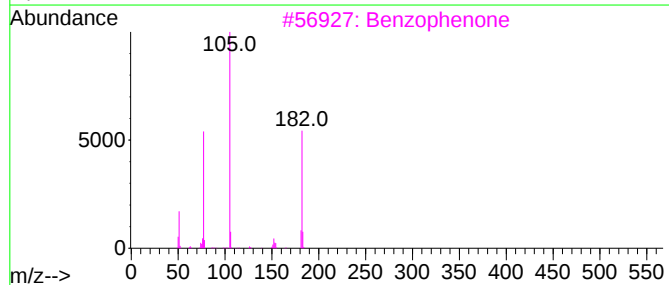
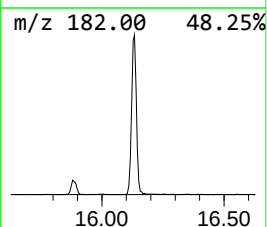
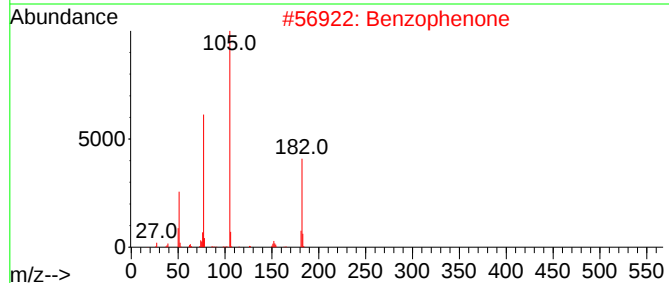
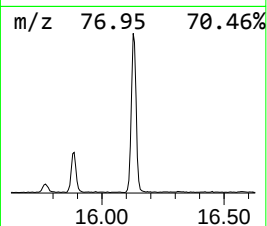
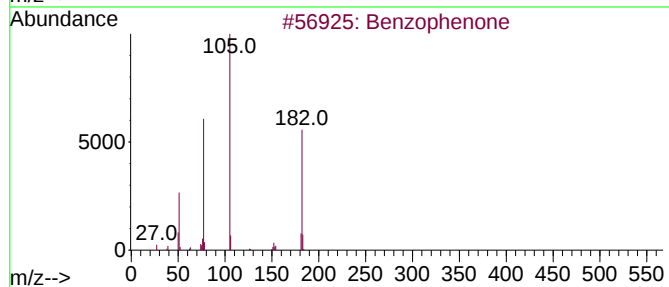
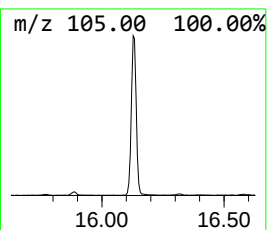
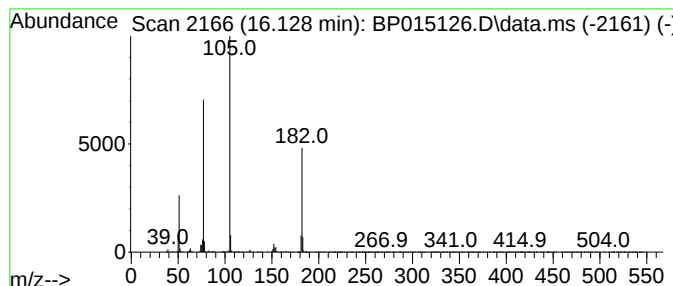
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051123.MA.M
Quant Title : SVOA CALIBRATION

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.
16.128	2.62 ng/ul	473982	Acenaphthene-d10	14.775

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	96
2			Benzophenone	182	C13H10O	000119-61-9	96
3			Benzophenone	182	C13H10O	000119-61-9	95
4			Benzophenone	182	C13H10O	000119-61-9	91
5			Benzophenone	182	C13H10O	000119-61-9	91



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzophenone	16.128	2.6	ng/u1	473982	3	14.775	3622670	20.0