

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051823\
 Data File : BP015146.D
 Acq On : 18 May 2023 15:34
 Operator : CG/JU
 Sample : 02612-22
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 JREJ3

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: BP015146.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	17	rVB	49367	73506	0.91%	0.084%
2	3.899	86	87	97	rBV	35928	63579	0.79%	0.073%
3	4.116	118	124	131	rBV	36399	63395	0.79%	0.073%
4	4.216	137	141	145	rVB	18448	24741	0.31%	0.028%
5	4.605	202	207	213	rVV4	12583	22732	0.28%	0.026%
6	4.675	215	219	226	rVB	50344	70469	0.87%	0.081%
7	5.169	298	303	309	rVB	25292	41905	0.52%	0.048%
8	5.293	320	324	328	rVB6	6831	10698	0.13%	0.012%
9	5.663	382	387	391	rBV	14743	22872	0.28%	0.026%
10	5.787	402	408	413	rBV	26062	40899	0.51%	0.047%
11	6.081	454	458	463	rVB6	5113	8301	0.10%	0.010%
12	7.234	648	654	656	rBV	38261	58526	0.73%	0.067%
13	7.275	656	661	674	rVB	1148114	1865395	23.14%	2.139%
14	7.446	681	690	702	rBV	995731	1581826	19.62%	1.814%
15	7.652	718	725	737	rBV	1150948	1853739	22.99%	2.126%
16	7.734	737	739	748	rVB8	9416	17571	0.22%	0.020%
17	7.852	754	759	765	rVB6	7052	11369	0.14%	0.013%
18	7.940	770	774	781	rVB4	7687	13423	0.17%	0.015%
19	8.128	799	806	814	rBV	1135655	1809437	22.44%	2.075%
20	8.840	918	927	941	rBV	1338764	2283937	28.33%	2.619%
21	8.963	942	948	955	rVB	196713	316031	3.92%	0.362%
22	9.304	998	1006	1015	rBV	1153892	1871837	23.22%	2.146%
23	9.463	1029	1033	1038	rVB7	5960	9823	0.12%	0.011%
24	9.699	1066	1073	1080	rVB	118240	179688	2.23%	0.206%
25	10.034	1116	1130	1140	rBV	942373	1601723	19.87%	1.837%
26	10.199	1150	1158	1166	rBV	7501	24939	0.31%	0.029%
27	10.393	1185	1191	1195	rBV2	16840	32719	0.41%	0.038%
28	10.428	1195	1197	1203	rVB6	9179	12675	0.16%	0.015%
29	10.569	1214	1221	1234	rBV	1422302	2433910	30.19%	2.791%
30	10.957	1277	1287	1296	rBV	1879515	3083873	38.25%	3.536%
31	11.093	1302	1310	1325	rBV	1371771	2432146	30.17%	2.789%
32	11.504	1374	1380	1388	rBV	6746	17630	0.22%	0.020%
33	11.763	1413	1424	1434	rBV10	10465	36856	0.46%	0.042%
34	12.193	1491	1497	1506	rVB	73372	110126	1.37%	0.126%
35	12.357	1520	1525	1530	rVB9	5294	10729	0.13%	0.012%
36	12.445	1535	1540	1543	rBV4	6034	11871	0.15%	0.014%

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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

37	12.534	1550	1555	1563	rVB2	19581	33881	0.42%	0.039%
38	13.581	1726	1733	1738	rBV3	14325	25366	0.31%	0.029%
39	13.651	1740	1745	1752	rVB4	11677	21604	0.27%	0.025%
40	13.739	1752	1760	1764	rBV7	6160	14795	0.18%	0.017%
41	14.087	1815	1819	1824	rVB8	10202	16141	0.20%	0.019%
42	14.169	1824	1833	1848	rBV	3251682	4325676	53.65%	4.960%
43	14.287	1848	1853	1858	rVB8	10383	17287	0.21%	0.020%
44	14.463	1875	1883	1898	rBV	3354625	4886314	60.61%	5.603%
45	14.645	1910	1914	1918	rVB5	5818	9435	0.12%	0.011%
46	14.769	1928	1935	1944	rBV	2942699	4347416	53.92%	4.985%
47	14.951	1958	1966	1986	rBV	1195008	1956265	24.26%	2.243%
48	15.175	1998	2004	2013	rVB2	49602	87442	1.08%	0.100%
49	15.716	2086	2096	2097	rBV	57985	71174	0.88%	0.082%
50	15.757	2097	2103	2114	rVV	4520232	6305148	78.21%	7.230%
51	15.869	2116	2122	2134	rBV	1223661	1730149	21.46%	1.984%
52	15.998	2139	2144	2148	rVB2	18829	31410	0.39%	0.036%
53	16.122	2157	2165	2173	rBV	871153	1208744	14.99%	1.386%
54	16.457	2219	2222	2228	rVB8	5909	10054	0.12%	0.012%
55	16.686	2257	2261	2266	rVB3	11641	16508	0.20%	0.019%
56	16.904	2294	2298	2302	rBV7	6670	13606	0.17%	0.016%
57	16.969	2305	2309	2313	rVV2	49808	65652	0.81%	0.075%
58	17.045	2318	2322	2327	rVB3	21598	32841	0.41%	0.038%
59	17.192	2342	2347	2352	rVB8	8209	13298	0.16%	0.015%
60	17.257	2355	2358	2361	rBV5	5897	8606	0.11%	0.010%
61	17.304	2364	2366	2372	rVB2	8532	12450	0.15%	0.014%
62	17.522	2396	2403	2413	rBV	3552508	5137569	63.72%	5.891%
63	17.616	2413	2419	2430	rVV2	4796136	6947491	86.17%	7.967%
64	17.704	2430	2434	2445	rVB4	86312	140993	1.75%	0.162%
65	17.857	2457	2460	2463	rBV5	6140	8848	0.11%	0.010%
66	17.892	2463	2466	2473	rVB8	6722	11074	0.14%	0.013%
67	18.004	2480	2485	2486	rBV5	9835	13128	0.16%	0.015%
68	18.033	2487	2490	2497	rVB2	33688	45237	0.56%	0.052%
69	18.163	2509	2512	2517	rVB7	9126	11500	0.14%	0.013%
70	18.269	2525	2530	2537	rVB6	21679	36628	0.45%	0.042%
71	18.375	2543	2548	2556	rBV	339937	477108	5.92%	0.547%
72	18.939	2640	2644	2648	rVB	60888	70478	0.87%	0.081%
73	19.151	2677	2680	2685	rBV2	27218	33814	0.42%	0.039%
74	19.416	2721	2725	2731	rVB3	73838	102848	1.28%	0.118%
75	19.480	2733	2736	2745	rVB2	61342	91020	1.13%	0.104%
76	19.598	2752	2756	2761	rBV	149159	199274	2.47%	0.229%

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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

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Peak separation: 5

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

77	19.704	2771	2774	2778	rVB2	32422	38654	0.48%	0.044%
78	19.839	2791	2797	2806	rBV	5809056	8062247	100.00%	9.245%
79	21.280	3037	3042	3051	rBV3	216769	444336	5.51%	0.510%
80	21.592	3090	3095	3110	rVB	3872675	5103825	63.31%	5.853%
81	22.210	3197	3200	3204	rVB	128136	137689	1.71%	0.158%
82	23.333	3388	3391	3398	rVB	219880	315280	3.91%	0.362%
83	23.992	3495	3503	3517	rBV2	3508413	7498221	93.00%	8.598%
84	24.157	3524	3531	3543	rVB	2300199	4891157	60.67%	5.609%

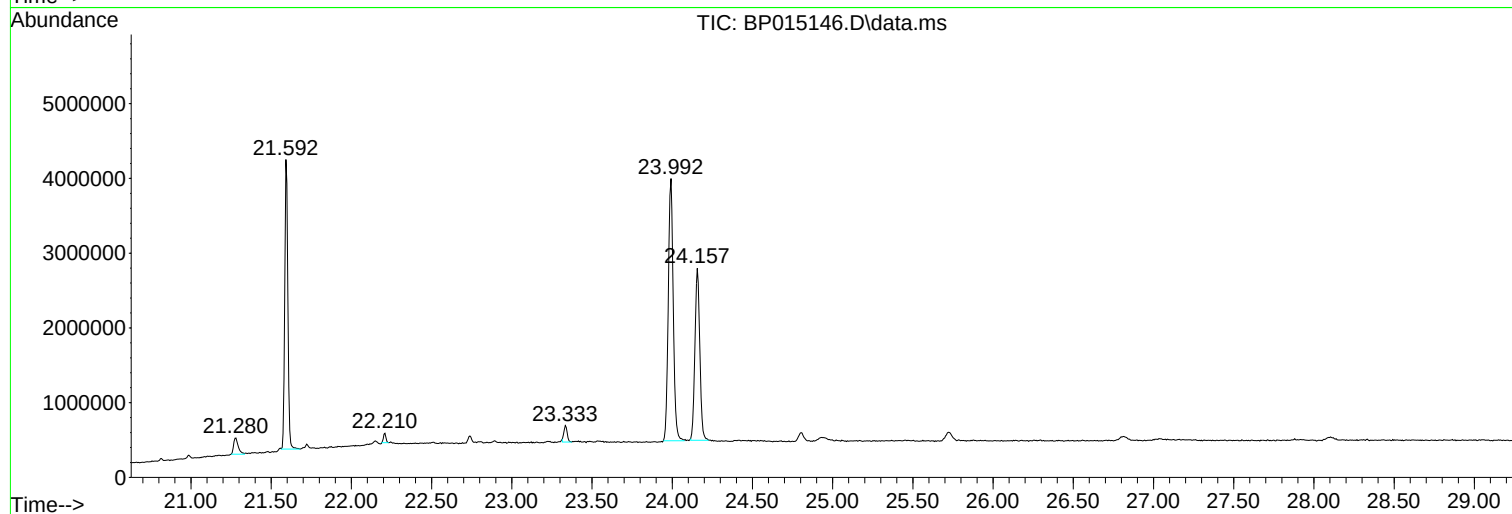
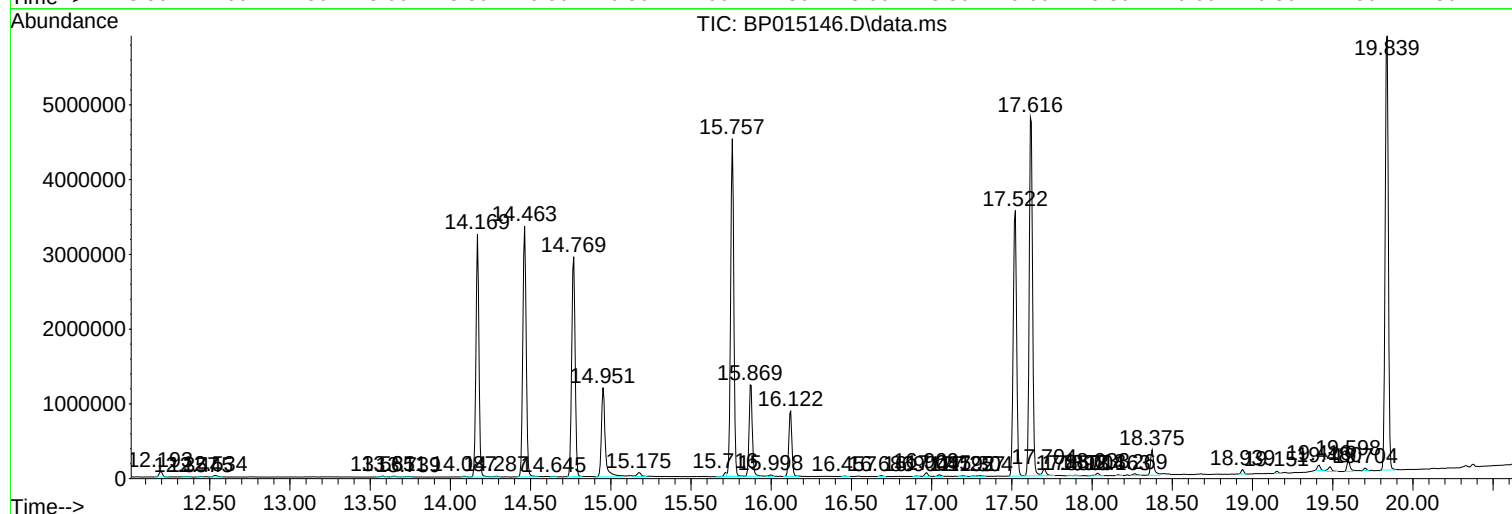
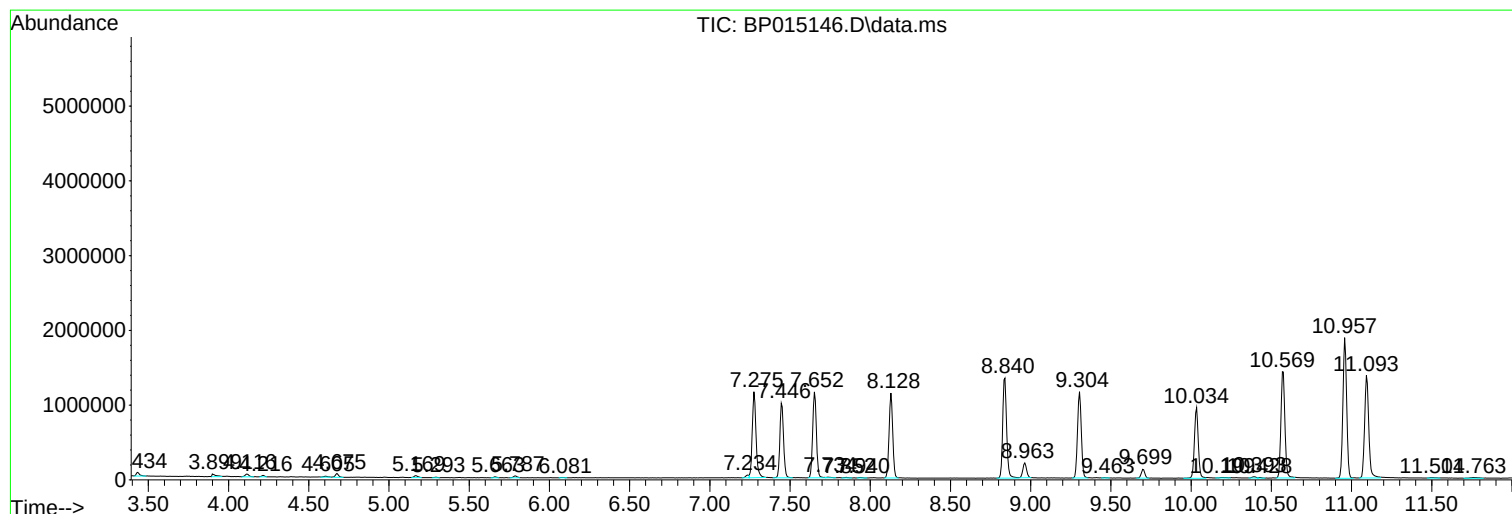
Sum of corrected areas: 87204577

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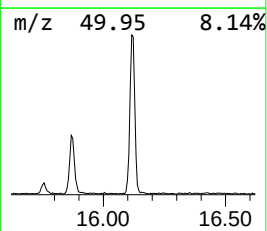
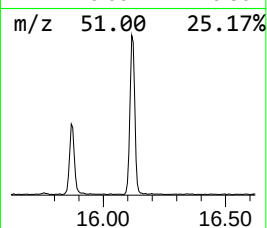
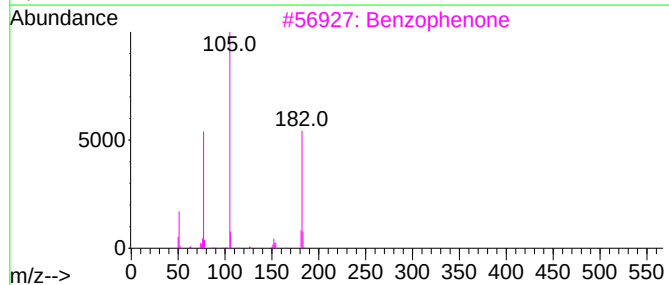
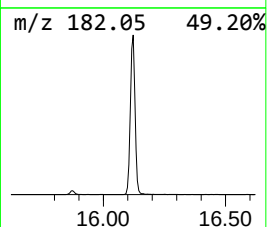
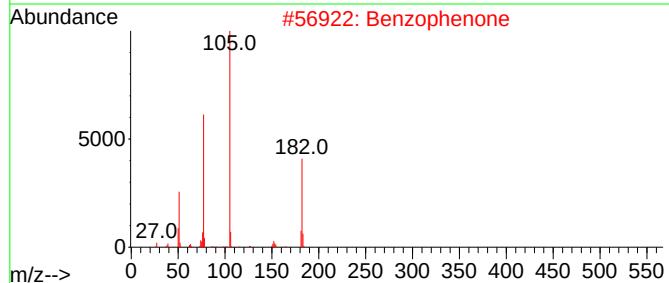
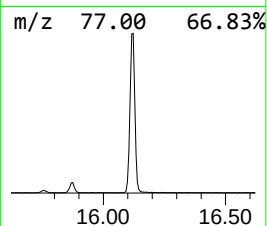
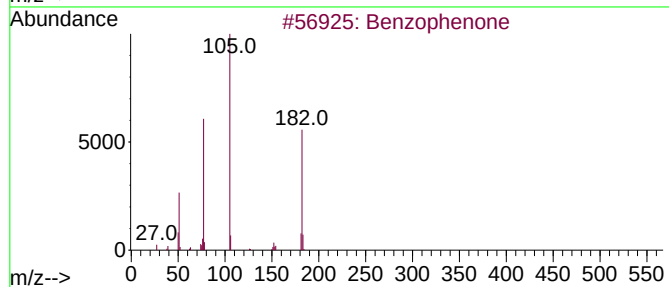
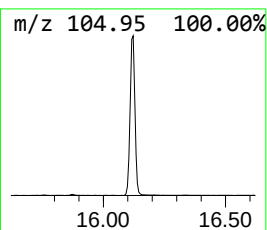
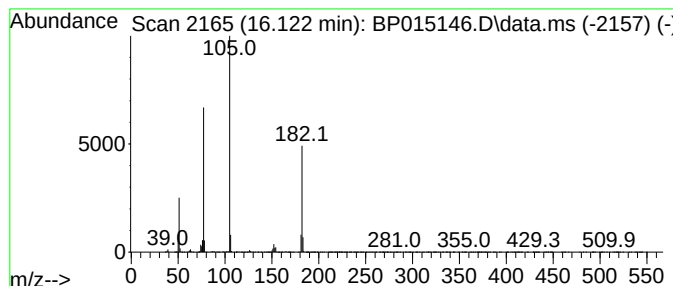
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051123.MA.M
Quant Title : SVOA CALIBRATION

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.
16.122	5.56 ng/ul	1208740	Acenaphthene-d10	14.769

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	90



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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	16.122	5.6	ng/ul	1208740	3	14.769	4347420	20.0