

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP060823\
 Data File : BP015433.D
 Acq On : 08 Jun 2023 13:15
 Operator : MA/JU
 Sample : 02960-04
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 H0FN8

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-BP060723.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP015433.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.411	35	38	44	rVB	15836	21598	0.82%	0.070%
2	3.905	119	122	123	rBV3	7253	6901	0.26%	0.022%
3	3.964	128	132	141	rVB	13597	23991	0.92%	0.077%
4	4.081	150	152	159	rVB3	9521	15277	0.58%	0.049%
5	4.187	168	170	173	rVB4	6842	7513	0.29%	0.024%
6	5.134	330	331	334	rBV3	3253	3351	0.13%	0.011%
7	5.264	349	353	355	rBV5	2587	3233	0.12%	0.010%
8	6.993	646	647	652	rBV5	1839	2678	0.10%	0.009%
9	7.252	684	691	708	rBV	233677	446741	17.04%	1.441%
10	7.416	714	719	729	rVB	197351	327443	12.49%	1.056%
11	7.622	747	754	763	rBV	260662	427495	16.31%	1.379%
12	8.099	827	835	845	rBV	456952	734100	28.01%	2.367%
13	8.252	860	861	867	rBV6	1772	3113	0.12%	0.010%
14	8.516	903	906	909	rVB5	2152	2909	0.11%	0.009%
15	8.810	949	956	971	rBV	219168	421703	16.09%	1.360%
16	8.934	971	977	985	rVV	90325	158204	6.04%	0.510%
17	9.275	1028	1035	1046	rBV	264941	460152	17.56%	1.484%
18	9.381	1051	1053	1057	rVB5	2668	3301	0.13%	0.011%
19	9.657	1096	1100	1106	rVB3	11020	16027	0.61%	0.052%
20	10.005	1150	1159	1179	rBV	229784	428525	16.35%	1.382%
21	10.246	1193	1200	1203	rBV9	2195	4546	0.17%	0.015%
22	10.369	1214	1221	1229	rVB5	9374	19992	0.76%	0.064%
23	10.481	1238	1240	1245	rBV6	2172	3333	0.13%	0.011%
24	10.552	1245	1252	1266	rBV	276625	568631	21.69%	1.834%
25	10.922	1309	1315	1325	rBV	663666	1132977	43.23%	3.654%
26	11.069	1333	1340	1356	rBV	203850	495355	18.90%	1.597%
27	11.381	1388	1393	1395	rBV6	2098	2826	0.11%	0.009%
28	11.504	1405	1414	1417	rBV10	3692	10839	0.41%	0.035%
29	11.769	1456	1459	1461	rBV4	2279	2803	0.11%	0.009%
30	12.157	1519	1525	1529	rBV5	6951	11185	0.43%	0.036%
31	12.328	1551	1554	1560	rBV7	2651	4510	0.17%	0.015%
32	12.510	1582	1585	1591	rBV6	5273	9018	0.34%	0.029%
33	12.657	1605	1610	1612	rBV6	1996	3031	0.12%	0.010%
34	13.093	1680	1684	1686	rBV5	2053	2945	0.11%	0.009%
35	13.546	1758	1761	1764	rBV5	2923	4038	0.15%	0.013%
36	13.622	1769	1774	1780	rVB5	8786	13676	0.52%	0.044%

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Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
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37	13.704	1783	1788	1789	rBV5	2634	3883	0.15%	0.013%
38	14.075	1848	1851	1852	rBV3	2578	2957	0.11%	0.010%
39	14.145	1855	1863	1878	rVV	821556	1189441	45.38%	3.836%
40	14.281	1878	1886	1898	rVV	36518	94577	3.61%	0.305%
41	14.434	1906	1912	1924	rBV2	822805	1242489	47.40%	4.007%
42	14.740	1957	1964	1976	rBV	1152479	1733813	66.15%	5.591%
43	14.957	1994	2001	2025	rBV2	91017	331457	12.65%	1.069%
44	15.169	2034	2037	2042	rVB7	9981	17769	0.68%	0.057%
45	15.234	2047	2048	2056	rVB8	3289	5015	0.19%	0.016%
46	15.428	2078	2081	2085	rVB5	2226	2773	0.11%	0.009%
47	15.563	2103	2104	2108	rVB4	3484	2648	0.10%	0.009%
48	15.681	2120	2124	2126	rBV4	5698	6616	0.25%	0.021%
49	15.734	2126	2133	2147	rVB	1136359	1660968	63.37%	5.356%
50	15.863	2148	2155	2164	rBV	238934	413836	15.79%	1.335%
51	15.975	2170	2174	2179	rVB6	8382	15339	0.59%	0.049%
52	16.098	2189	2195	2208	rBV	360144	498493	19.02%	1.608%
53	16.598	2277	2280	2281	rBV3	2397	2729	0.10%	0.009%
54	16.663	2287	2291	2295	rVB	13470	17335	0.66%	0.056%
55	16.939	2333	2338	2347	rVB5	7654	14257	0.54%	0.046%
56	17.028	2347	2353	2359	rBV	33830	53652	2.05%	0.173%
57	17.175	2374	2378	2381	rBV6	3989	6309	0.24%	0.020%
58	17.228	2385	2387	2392	rVV5	3368	6511	0.25%	0.021%
59	17.286	2392	2397	2402	rVB9	4947	11874	0.45%	0.038%
60	17.398	2411	2416	2419	rBV7	4461	8061	0.31%	0.026%
61	17.498	2426	2433	2443	rBV	1557046	2284175	87.15%	7.366%
62	17.598	2443	2450	2460	rVV2	1275146	1918892	73.21%	6.188%
63	17.769	2473	2479	2483	rBV4	6776	12144	0.46%	0.039%
64	17.998	2514	2518	2524	rVB6	6801	11392	0.43%	0.037%
65	18.245	2555	2560	2566	rBV6	13840	27120	1.03%	0.087%
66	18.351	2574	2578	2585	rBV	162699	223409	8.52%	0.720%
67	18.745	2643	2645	2647	rBV2	6249	5472	0.21%	0.018%
68	18.922	2671	2675	2679	rVB3	17989	26810	1.02%	0.086%
69	19.139	2708	2712	2716	rBV	20520	29020	1.11%	0.094%
70	19.181	2716	2719	2723	rVB5	14582	19838	0.76%	0.064%
71	19.392	2745	2755	2761	rBV8	48724	132130	5.04%	0.426%
72	19.469	2765	2768	2775	rVB3	25687	39589	1.51%	0.128%
73	19.580	2783	2787	2794	rBV2	51839	83041	3.17%	0.268%
74	19.816	2821	2827	2838	rBV	1909313	2575169	98.25%	8.304%
75	20.757	2983	2987	2990	rBV2	74716	88631	3.38%	0.286%
76	21.251	3067	3071	3078	rBV3	139320	258328	9.86%	0.833%

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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

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Title : SVOA CALIBRATION

77	21.457	3101	3106	3114	rBV	2205102	2566356	97.91%	8.276%
78	21.580	3121	3127	3135	rBV	1822029	2621083	100.00%	8.452%
79	23.963	3525	3532	3543	rVB2	1155345	2447377	93.37%	7.892%
80	24.133	3554	3561	3569	rVB	1134193	2493101	95.12%	8.040%

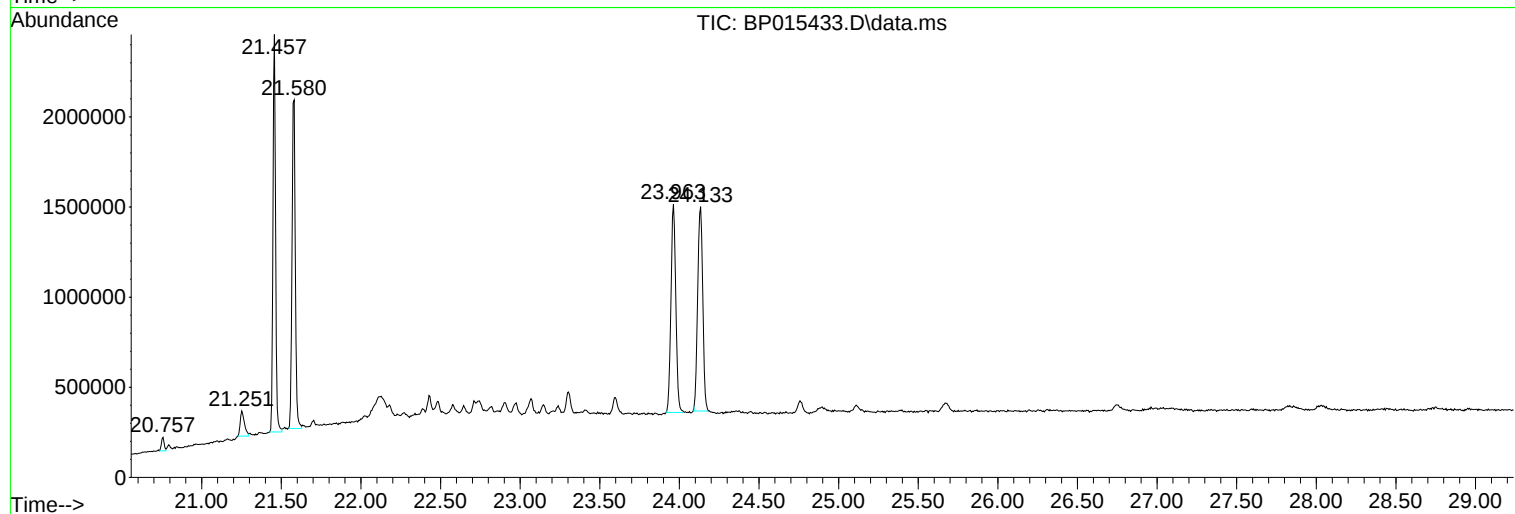
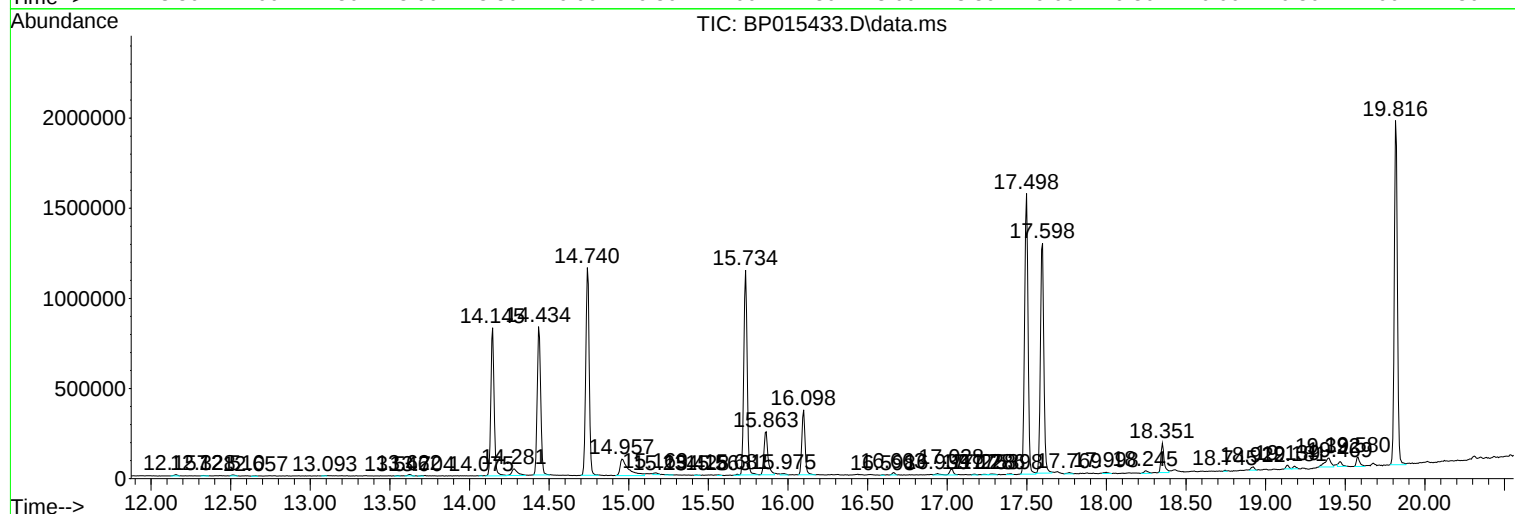
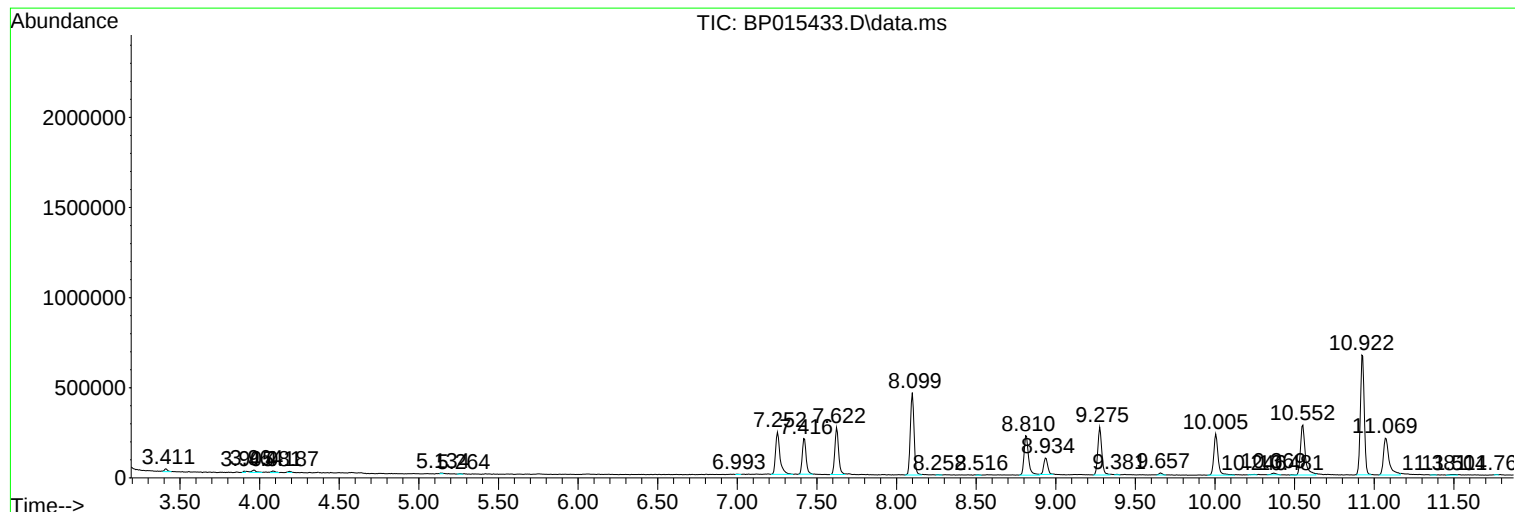
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

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



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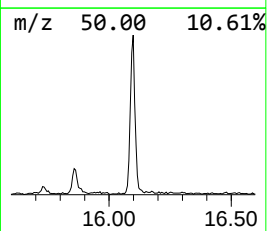
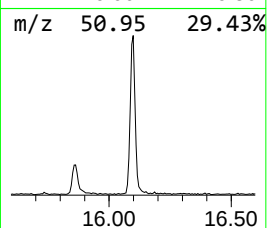
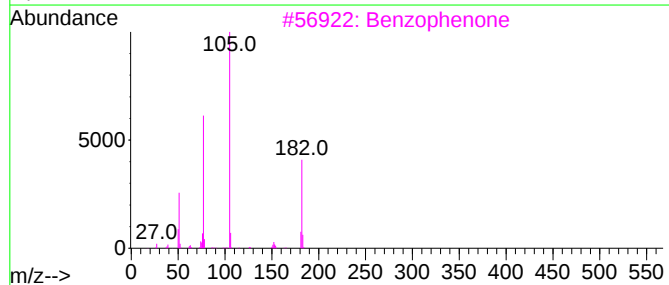
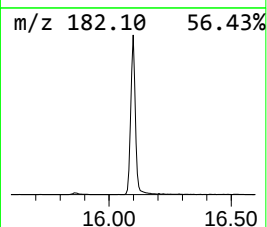
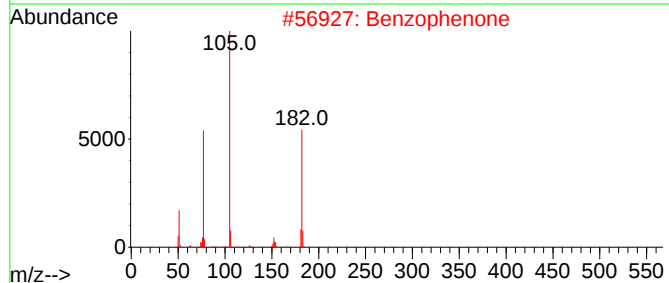
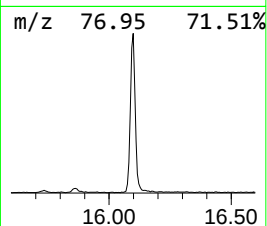
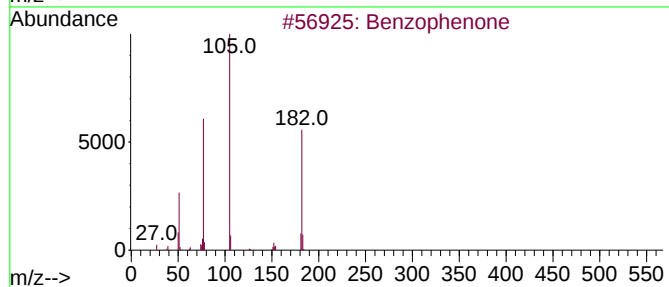
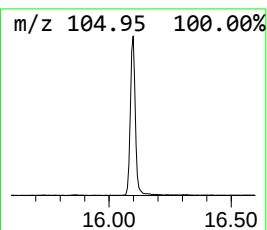
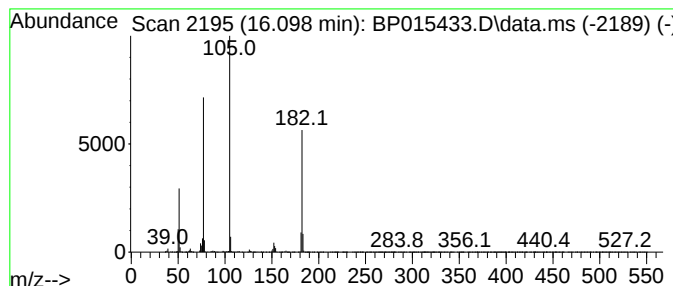
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

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.
16.098	5.75 ng/ul	498493	Acenaphthene-d10	14.740

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	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	16.098	5.8	ng/u1	498493	3	14.740	1733810	20.0