

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052323\
 Data File : BP015239.D
 Acq On : 23 May 2023 14:18
 Operator : CG/JU
 Sample : 02861-06
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 H0FL1

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: BP015239.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.428	37	41	48	rVB	38262	51361	1.87%	0.180%
2	3.881	117	118	133	rBV	71103	133268	4.86%	0.467%
3	4.105	153	156	161	rVB5	5555	10544	0.38%	0.037%
4	4.211	170	174	180	rVB2	12713	18857	0.69%	0.066%
5	4.663	247	251	258	rVB4	5741	9833	0.36%	0.034%
6	5.658	418	420	423	rVB4	2965	3037	0.11%	0.011%
7	5.775	438	440	447	rVB5	3653	4510	0.16%	0.016%
8	6.552	568	572	575	rBV6	2628	4388	0.16%	0.015%
9	7.222	682	686	689	rVB5	4589	6401	0.23%	0.022%
10	7.269	689	694	705	rVB	50568	96564	3.52%	0.338%
11	7.440	716	723	737	rVB	298163	472503	17.24%	1.654%
12	7.646	751	758	768	rBV	265093	436150	15.92%	1.527%
13	8.122	832	839	852	rBV	729651	1172895	42.80%	4.107%
14	8.834	953	960	972	rBV	175871	328989	12.01%	1.152%
15	9.299	1032	1039	1052	rBV	354350	620996	22.66%	2.174%
16	9.446	1062	1064	1065	rBV2	3603	2939	0.11%	0.010%
17	9.687	1100	1105	1113	rVB	15137	24579	0.90%	0.086%
18	10.022	1155	1162	1172	rVV	265818	465273	16.98%	1.629%
19	10.563	1245	1254	1275	rBV	378894	716271	26.14%	2.508%
20	10.951	1310	1320	1329	rBV	1042933	1783001	65.07%	6.243%
21	11.087	1334	1343	1360	rBV	252897	502020	18.32%	1.758%
22	11.393	1394	1395	1403	rVB8	2111	3320	0.12%	0.012%
23	12.181	1523	1529	1535	rVB4	13330	20099	0.73%	0.070%
24	12.540	1582	1590	1595	rBV5	5844	11115	0.41%	0.039%
25	12.751	1623	1626	1629	rBV5	3558	4626	0.17%	0.016%
26	13.363	1726	1730	1733	rVV6	2468	3749	0.14%	0.013%
27	13.569	1759	1765	1768	rVB8	3583	5473	0.20%	0.019%
28	13.645	1772	1778	1784	rBV	22765	37128	1.35%	0.130%
29	13.728	1788	1792	1795	rBV6	1794	3150	0.11%	0.011%
30	14.163	1858	1866	1876	rBV	907487	1331624	48.59%	4.663%
31	14.275	1881	1885	1892	rVV9	4588	10347	0.38%	0.036%
32	14.339	1892	1896	1901	rVB3	4488	7003	0.26%	0.025%
33	14.451	1908	1915	1928	rBV	995716	1509062	55.07%	5.284%
34	14.757	1958	1967	1975	rBV	1655939	2460285	89.78%	8.614%
35	14.992	2005	2007	2014	rVB8	3568	6278	0.23%	0.022%
36	15.163	2033	2036	2040	rVV5	2354	3234	0.12%	0.011%

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

37	15.375	2069	2072	2077	rVV7	1928	3051	0.11%	0.011%
38	15.533	2095	2099	2101	rVV4	2333	3547	0.13%	0.012%
39	15.563	2101	2104	2106	rVV4	3639	4648	0.17%	0.016%
40	15.704	2123	2128	2130	rBV4	8805	12536	0.46%	0.044%
41	15.745	2130	2135	2147	rVV	1244641	1905076	69.52%	6.670%
42	15.869	2150	2156	2164	rBV	86869	145089	5.29%	0.508%
43	15.992	2172	2177	2182	rVB6	7991	13121	0.48%	0.046%
44	16.110	2191	2197	2210	rBV	315558	449152	16.39%	1.573%
45	16.445	2252	2254	2259	rBV6	3657	4884	0.18%	0.017%
46	16.675	2289	2293	2299	rBV	37690	55617	2.03%	0.195%
47	16.904	2328	2332	2333	rBV4	3054	3885	0.14%	0.014%
48	16.963	2337	2342	2345	rVV5	6893	9696	0.35%	0.034%
49	17.086	2361	2363	2367	rVB4	2295	3058	0.11%	0.011%
50	17.122	2367	2369	2373	rBV5	2296	3810	0.14%	0.013%
51	17.245	2385	2390	2396	rBV8	7050	12208	0.45%	0.043%
52	17.298	2396	2399	2404	rVB6	6407	8855	0.32%	0.031%
53	17.510	2428	2435	2443	rBV	1953745	2740258	100.00%	9.595%
54	17.610	2446	2452	2465	rVB	1364519	1972564	71.98%	6.907%
55	17.780	2477	2481	2484	rBV6	4731	8297	0.30%	0.029%
56	17.980	2512	2515	2519	rBV6	6415	7677	0.28%	0.027%
57	18.151	2542	2544	2550	rVB7	4892	6526	0.24%	0.023%
58	18.363	2575	2580	2587	rBV3	60983	93982	3.43%	0.329%
59	18.769	2647	2649	2654	rBV6	5713	8713	0.32%	0.031%
60	19.404	2754	2757	2762	rVB5	19434	26862	0.98%	0.094%
61	19.474	2767	2769	2773	rVB2	15806	19612	0.72%	0.069%
62	19.592	2786	2789	2792	rBV5	17532	21370	0.78%	0.075%
63	19.827	2824	2829	2838	rBV	1637452	2122646	77.46%	7.432%
64	21.586	3123	3128	3136	rBV	1637810	2245998	81.96%	7.864%
65	23.974	3527	3534	3545	rBV2	896459	2035659	74.29%	7.128%
66	24.139	3555	3562	3573	rVB	1067937	2330597	85.05%	8.160%

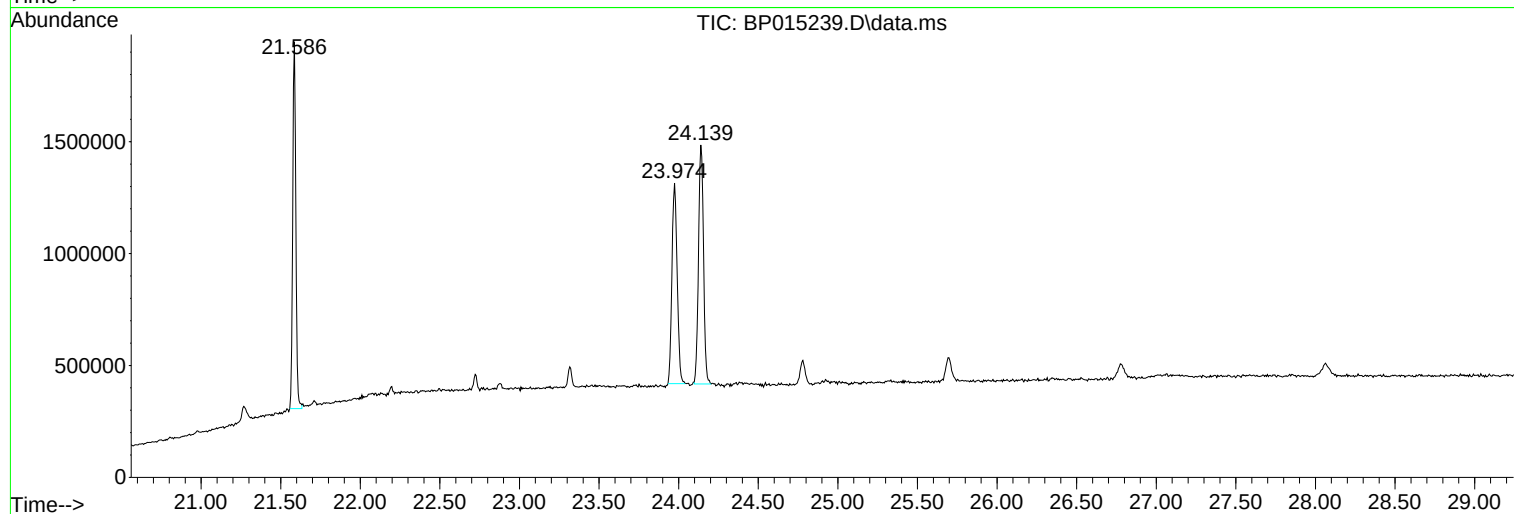
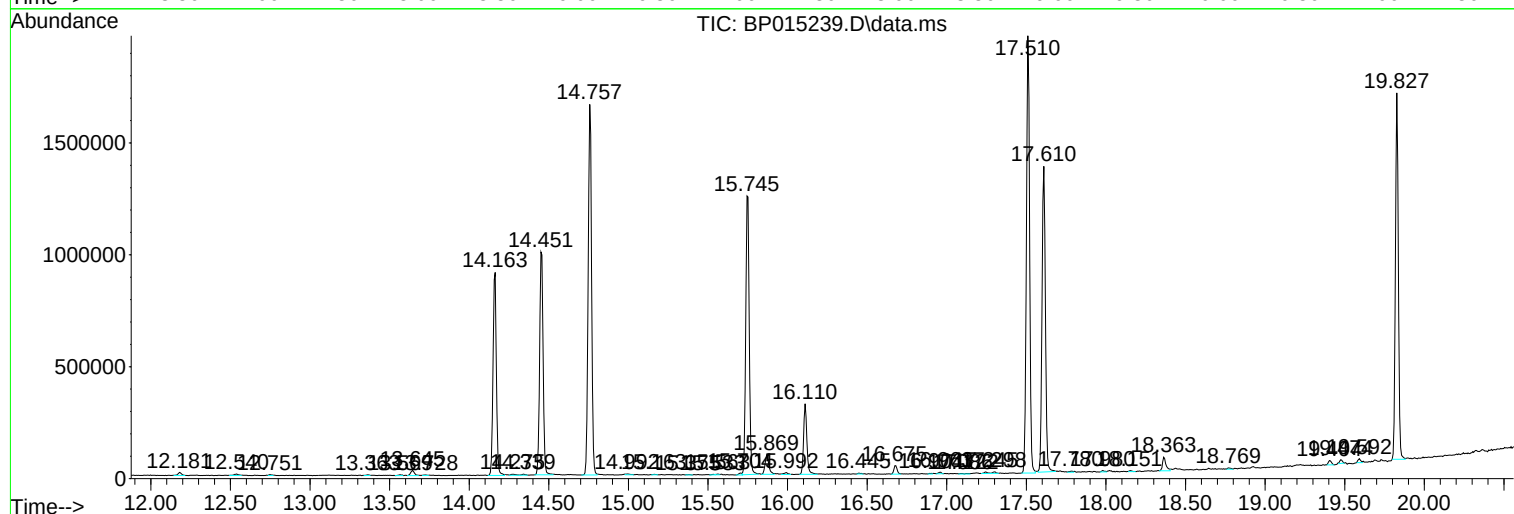
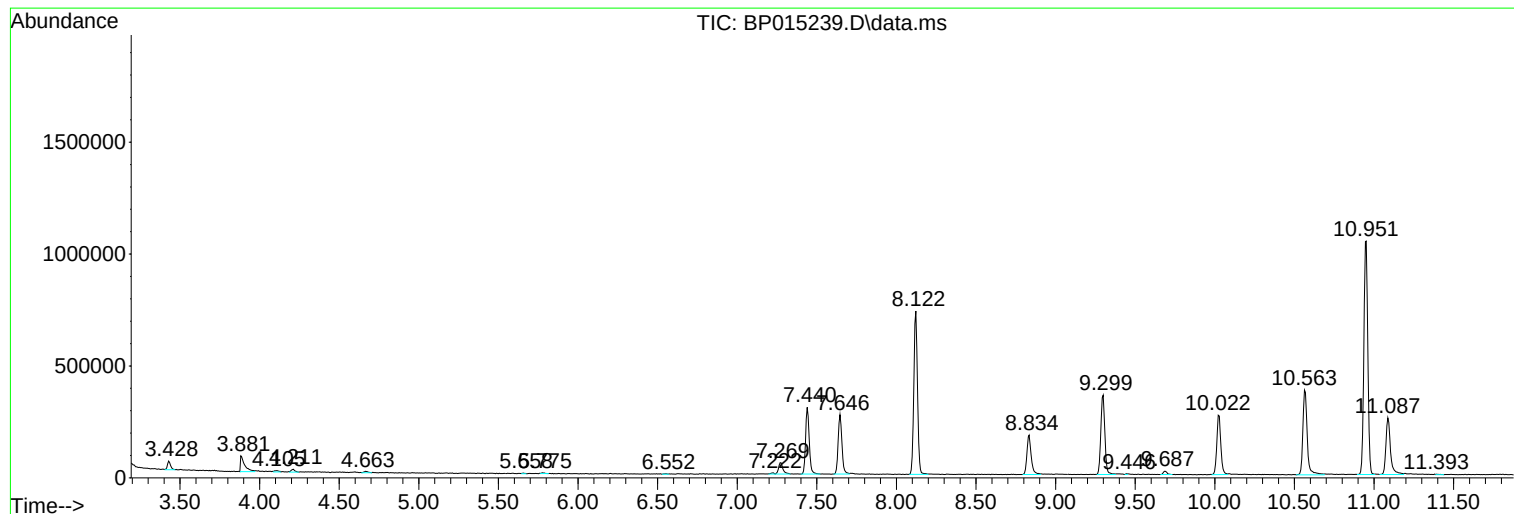
Sum of corrected areas: 28559866

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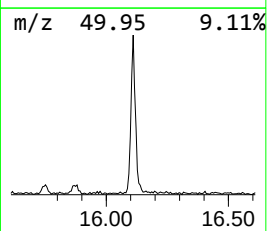
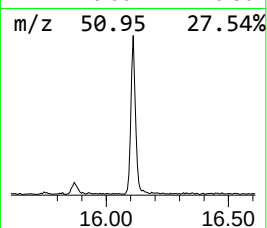
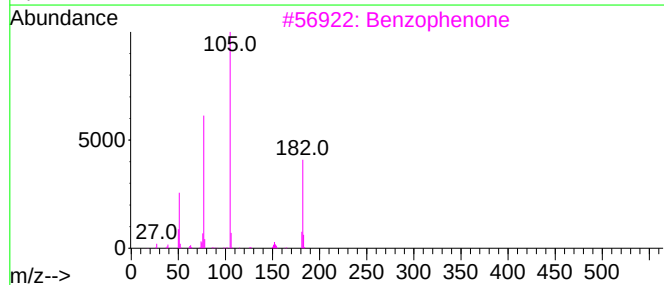
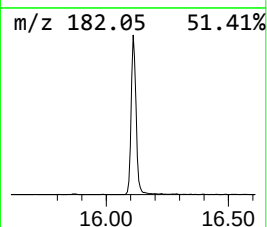
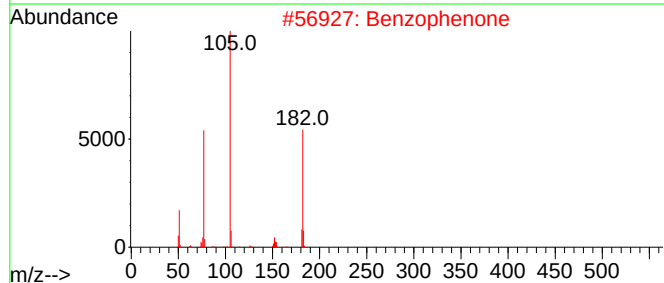
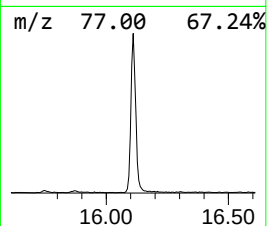
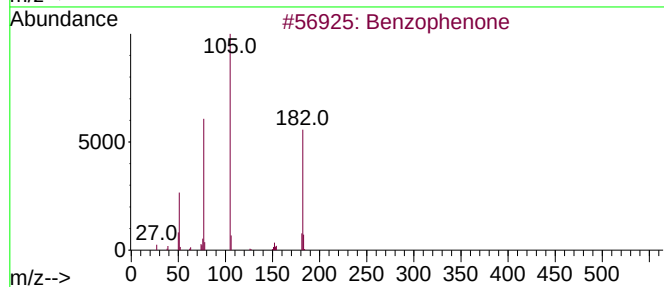
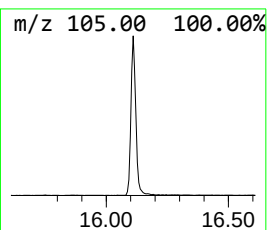
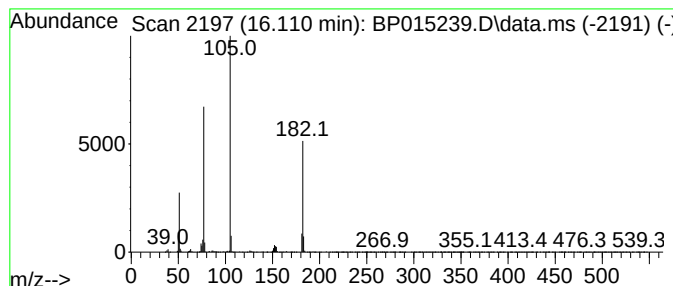
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051123.MA.M
Quant Title : SVOA CALIBRATION

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.
16.110	3.65 ng/ul	449152	Acenaphthene-d10	14.757

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	94
5			Benzophenone	182	C13H10O	000119-61-9	90



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzophenone	16.110	3.6	ng/u1	449152	3	14.757	2460290	20.0