

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111124\
 Data File : BM048593.D
 Acq On : 11 Nov 2024 20:21
 Operator : RC/JU
 Sample : P4714-10
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 COAH0

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_M\Methods\SFAM-EPA-BM110724.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BM048593.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.122	20	23	31	rVB	31595	45089	1.25%	0.137%
2	3.546	92	95	115	rBV	136890	263120	7.31%	0.798%
3	3.857	145	148	153	rBV3	6646	9277	0.26%	0.028%
4	4.369	231	235	243	rVB2	20123	28492	0.79%	0.086%
5	5.104	356	360	363	rBV5	3803	5663	0.16%	0.017%
6	5.310	390	395	403	rVB8	4690	12149	0.34%	0.037%
7	5.522	427	431	436	rBV4	6152	10069	0.28%	0.031%
8	5.645	445	452	454	rBV6	3373	7675	0.21%	0.023%
9	5.681	455	458	462	rVB6	5701	7878	0.22%	0.024%
10	5.775	471	474	477	rBV5	2589	4054	0.11%	0.012%
11	5.934	497	501	505	rBV5	2863	4648	0.13%	0.014%
12	6.222	546	550	553	rBV6	4075	6165	0.17%	0.019%
13	6.298	560	563	566	rVV4	3570	4239	0.12%	0.013%
14	6.322	566	567	573	rVB6	3234	4777	0.13%	0.014%
15	6.657	619	624	628	rVB8	3572	6431	0.18%	0.020%
16	6.845	643	656	672	rBV	108738	223831	6.22%	0.679%
17	6.998	675	682	694	rBV	365461	624934	17.36%	1.896%
18	7.192	709	715	726	rBV	418026	725965	20.17%	2.203%
19	7.281	727	730	737	rVB5	15176	26728	0.74%	0.081%
20	7.392	747	749	753	rVB5	4102	3929	0.11%	0.012%
21	7.657	787	794	804	rBV	508517	839661	23.32%	2.548%
22	7.945	841	843	849	rVB7	2612	3805	0.11%	0.012%
23	8.375	910	916	931	rBV	337121	628855	17.47%	1.908%
24	8.816	985	991	1000	rBV	419352	723118	20.09%	2.194%
25	9.410	1086	1092	1096	rBV7	2600	5361	0.15%	0.016%
26	9.539	1105	1114	1126	rBV	336033	615797	17.11%	1.869%
27	10.080	1199	1206	1228	rBV	484211	976744	27.13%	2.964%
28	10.445	1261	1268	1284	rBV	716325	1213580	33.71%	3.682%
29	10.592	1286	1293	1314	rBV	467253	1036702	28.80%	3.146%
30	11.157	1382	1389	1400	rVB6	15084	32456	0.90%	0.098%
31	12.069	1540	1544	1550	rVB5	7112	13725	0.38%	0.042%
32	12.821	1666	1672	1679	rBV8	5892	12199	0.34%	0.037%
33	13.168	1725	1731	1733	rVB6	3085	6110	0.17%	0.019%
34	13.721	1819	1825	1840	rVB	986075	1450791	40.30%	4.402%
35	13.998	1865	1872	1885	rBV	1151159	1707690	47.44%	5.182%
36	14.310	1918	1925	1932	rBV	1037385	1562051	43.39%	4.740%

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Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM110724.MA.M
Title : SVOA CALIBRATION

37	14.533	1958	1963	1965	rBV2	35885	50205	1.39%	0.152%
38	14.563	1965	1968	1981	rVB5	29494	85793	2.38%	0.260%
39	14.974	2034	2038	2041	rVB4	6488	9108	0.25%	0.028%
40	15.110	2058	2061	2063	rBV4	2712	3789	0.11%	0.011%
41	15.304	2086	2094	2102	rBV	1569024	2264106	62.89%	6.870%
42	15.439	2111	2117	2129	rBV	375476	598913	16.64%	1.817%
43	15.545	2131	2135	2143	rVB5	9717	19241	0.53%	0.058%
44	15.674	2151	2157	2166	rBV	194366	276748	7.69%	0.840%
45	15.862	2184	2189	2194	rBV5	7343	13557	0.38%	0.041%
46	16.051	2217	2221	2225	rVV3	10200	16208	0.45%	0.049%
47	16.233	2251	2252	2259	rVB6	5380	6650	0.18%	0.020%
48	16.474	2288	2293	2295	rBV6	2862	4556	0.13%	0.014%
49	16.545	2300	2305	2312	rBV7	10150	18239	0.51%	0.055%
50	16.845	2349	2356	2359	rBV6	4394	11459	0.32%	0.035%
51	16.880	2360	2362	2368	rVB5	4687	7400	0.21%	0.022%
52	17.051	2385	2391	2402	rBV	1274719	1869992	51.95%	5.674%
53	17.151	2402	2408	2421	rVV2	1823921	2733824	75.94%	8.295%
54	17.727	2503	2506	2509	rVB5	3765	4825	0.13%	0.015%
55	17.827	2516	2523	2531	rBV9	10815	26063	0.72%	0.079%
56	17.892	2531	2534	2537	rBV5	2449	4146	0.12%	0.013%
57	17.951	2539	2544	2550	rBV	80987	128813	3.58%	0.391%
58	18.245	2592	2594	2598	rVB5	5929	7925	0.22%	0.024%
59	19.015	2721	2725	2732	rVB3	26971	39267	1.09%	0.119%
60	19.086	2733	2737	2740	rBV	26226	37110	1.03%	0.113%
61	19.239	2759	2763	2767	rBV6	25679	37948	1.05%	0.115%
62	19.450	2793	2799	2811	rBV	2485665	3432920	95.36%	10.416%
63	20.950	3050	3054	3057	rBV2	106454	162638	4.52%	0.493%
64	21.280	3105	3110	3126	rVB	1376373	2146909	59.64%	6.514%
65	23.991	3562	3571	3585	rBV	1485073	3599921	100.00%	10.923%
66	24.191	3597	3605	3621	rVB	972802	2484552	69.02%	7.539%

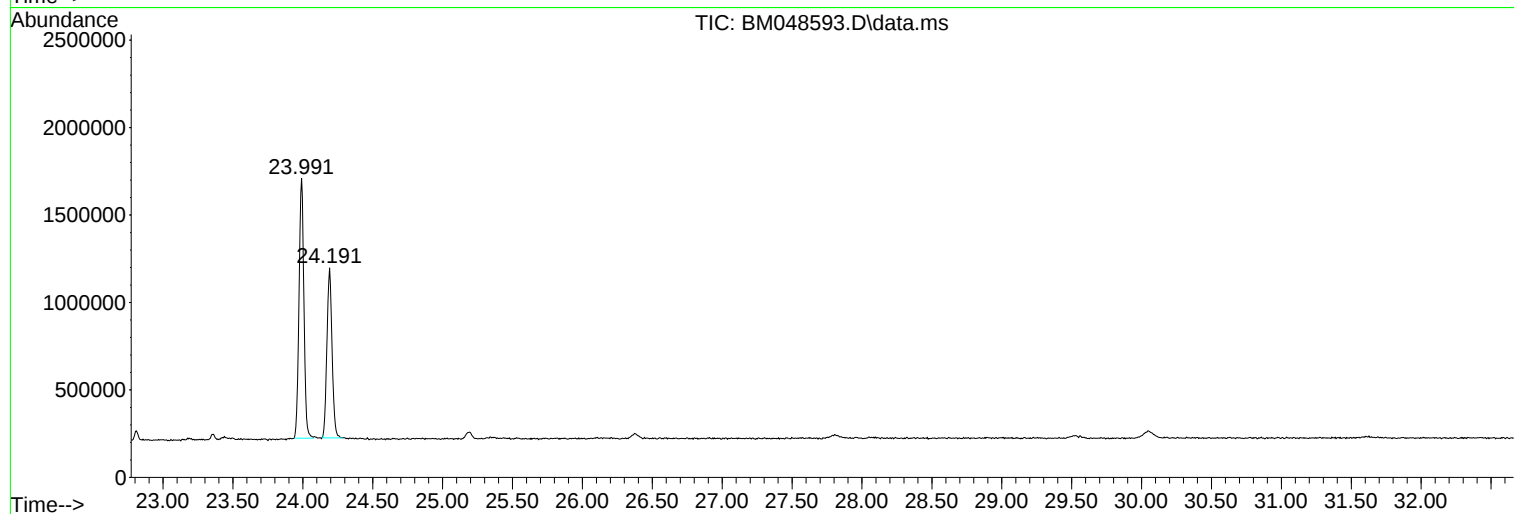
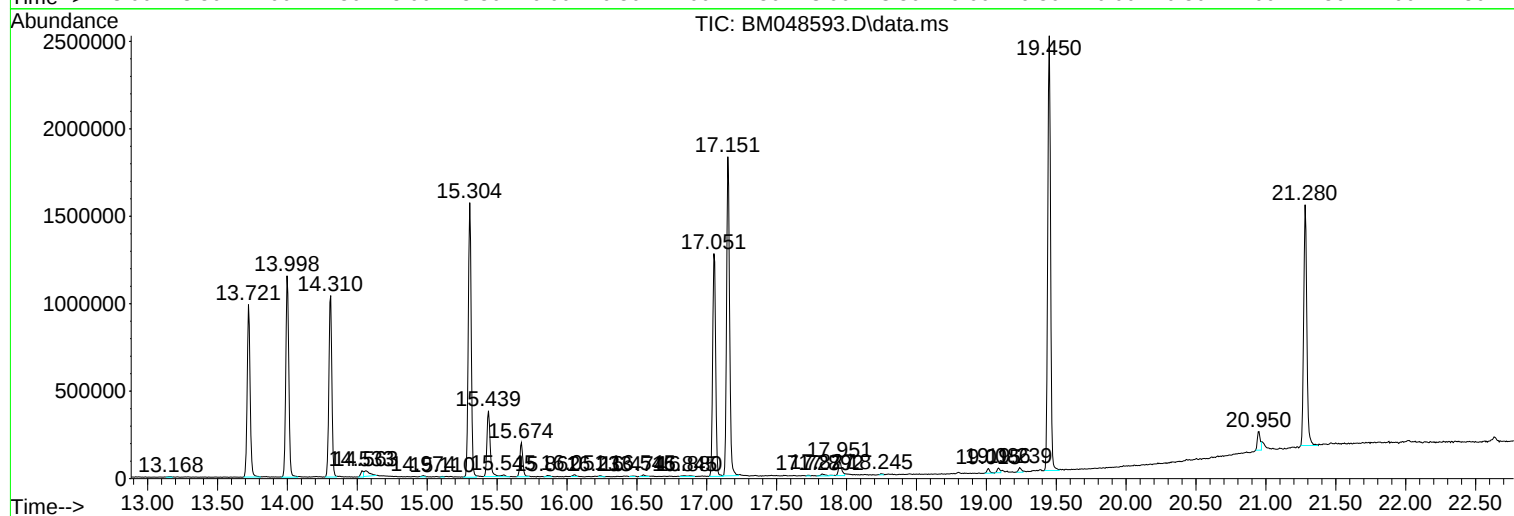
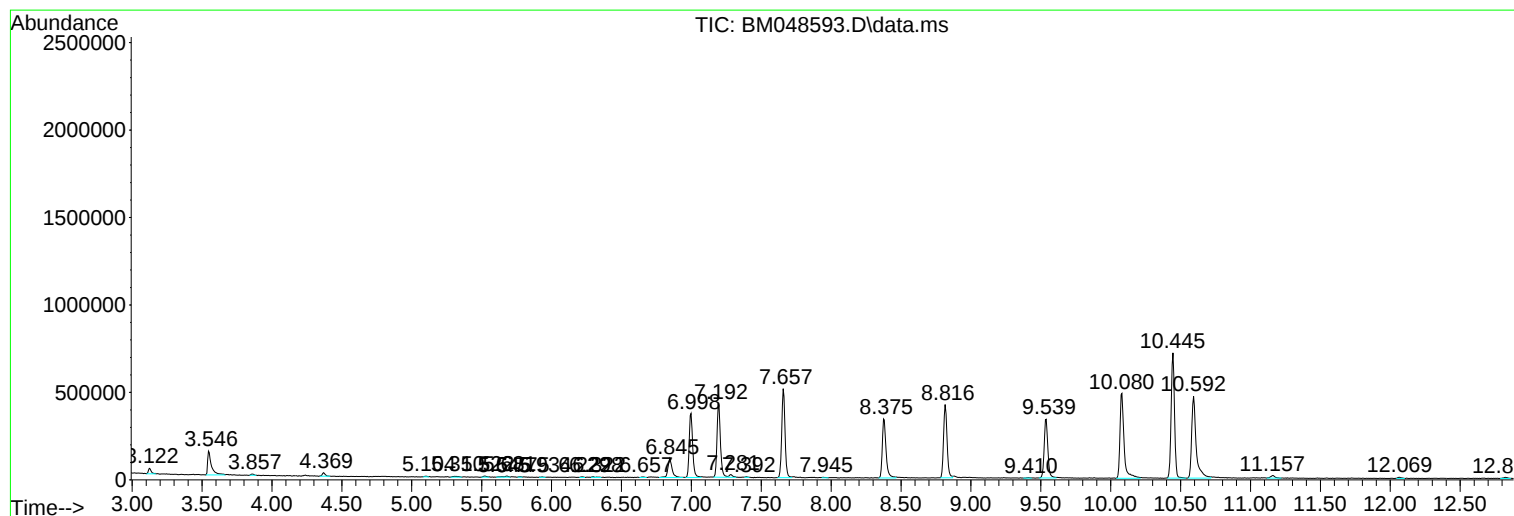
Sum of corrected areas: 32956583

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM110724.MA.M
Quant Title : SVOA CALIBRATION

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



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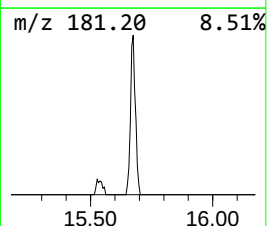
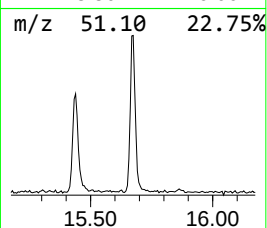
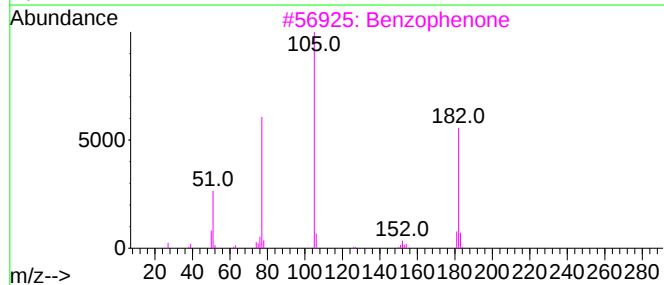
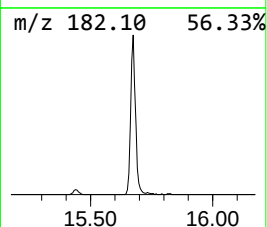
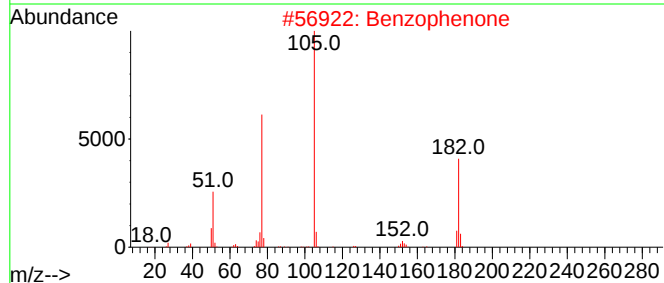
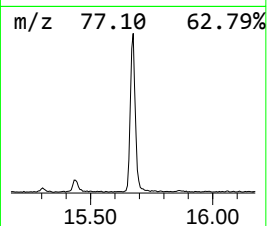
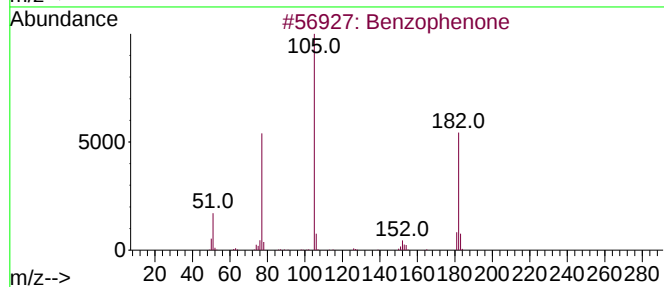
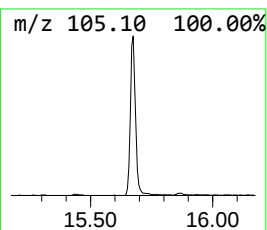
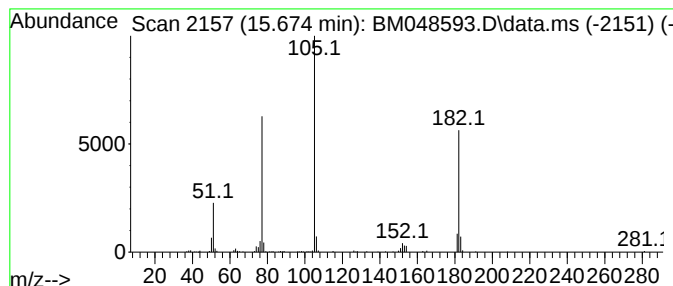
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM110724.MA.M
Quant Title : SVOA CALIBRATION

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.
15.674	3.54 ng/ul	276748	Acenaphthene-d10	14.310

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	96
4			Benzophenone	182	C13H10O	000119-61-9	91
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	15.674	3.5	ng/u1	276748	3	14.310	1562050	20.0