

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121024\
 Data File : BP023386.D
 Acq On : 11 Dec 2024 20:43
 Operator : RC/JU
 Sample : P5152-04
 Misc :
 ALS Vial : 50 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BHK73

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-BP112624.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP023386.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.099	16	19	29	rVB	21852	31473	1.07%	0.126%
2	3.528	88	92	104	rBV	55293	125803	4.27%	0.502%
3	3.822	138	142	147	rVB2	6512	9386	0.32%	0.037%
4	4.752	296	300	307	rBV2	7365	13010	0.44%	0.052%
5	6.675	622	627	633	rVB3	5857	9460	0.32%	0.038%
6	6.769	636	643	658	rVV	37111	123226	4.18%	0.492%
7	6.893	658	664	686	rVB	237322	390049	13.24%	1.556%
8	7.093	691	698	718	rBV	230360	447497	15.19%	1.785%
9	7.540	768	774	786	rBV	348822	595617	20.21%	2.376%
10	8.263	891	897	918	rBV	160751	357135	12.12%	1.425%
11	8.681	962	968	986	rBV	294382	527974	17.92%	2.106%
12	9.063	1026	1033	1042	rVB3	10909	21202	0.72%	0.085%
13	9.287	1067	1071	1076	rBV7	3780	7706	0.26%	0.031%
14	9.393	1083	1089	1106	rBV	205603	447539	15.19%	1.785%
15	9.945	1175	1183	1205	rBV	244882	639716	21.71%	2.552%
16	10.287	1233	1241	1259	rVB	475656	791281	26.85%	3.157%
17	10.445	1261	1268	1293	rBV	200716	552700	18.76%	2.205%
18	11.534	1445	1453	1458	rBV7	3141	6871	0.23%	0.027%
19	11.934	1514	1521	1529	rBV7	3003	7565	0.26%	0.030%
20	13.498	1781	1787	1790	rBV8	1620	3239	0.11%	0.013%
21	13.557	1792	1797	1818	rVV	819735	1154752	39.19%	4.607%
22	13.845	1839	1846	1867	rBV	874517	1162573	39.45%	4.638%
23	14.157	1891	1899	1912	rBV	730714	1171093	39.74%	4.672%
24	14.510	1953	1959	1976	rBV5	16763	83199	2.82%	0.332%
25	14.963	2032	2036	2040	rVB5	2160	3657	0.12%	0.015%
26	15.004	2040	2043	2047	rVB6	1896	3033	0.10%	0.012%
27	15.169	2063	2071	2085	rBV	868755	1626589	55.20%	6.489%
28	15.327	2092	2098	2111	rBV	280120	553575	18.79%	2.208%
29	15.416	2111	2113	2119	rVB6	10441	17106	0.58%	0.068%
30	15.545	2130	2135	2150	rVB	175389	265691	9.02%	1.060%
31	15.757	2165	2171	2176	rBV10	2205	4929	0.17%	0.020%
32	16.127	2227	2234	2239	rVB10	2139	4059	0.14%	0.016%
33	16.757	2336	2341	2345	rVB7	2775	4924	0.17%	0.020%
34	16.898	2362	2365	2370	rVB7	2492	3553	0.12%	0.014%
35	16.969	2370	2377	2387	rBV	1103655	1617223	54.88%	6.452%
36	17.063	2387	2393	2411	rVV	1343551	2098549	71.22%	8.372%

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37	17.204	2413	2417	2426	rVB6	13726	25589	0.87%	0.102%
38	17.910	2531	2537	2547	rBV	90995	153824	5.22%	0.614%
39	18.363	2606	2614	2616	rBV9	3828	8093	0.27%	0.032%
40	18.768	2681	2683	2686	rVB4	5315	5546	0.19%	0.022%
41	18.998	2717	2722	2727	rBV2	19385	27601	0.94%	0.110%
42	19.092	2733	2738	2742	rBV	15086	28326	0.96%	0.113%
43	19.239	2759	2763	2771	rBV3	32484	51561	1.75%	0.206%
44	19.392	2786	2789	2792	rBV5	7068	7367	0.25%	0.029%
45	19.445	2793	2798	2817	rBV	2095223	2773271	94.12%	11.064%
46	21.392	3123	3129	3146	rVB	1109730	2039490	69.22%	8.137%
47	24.351	3623	3632	3645	rBV	1161682	2946586	100.00%	11.755%
48	24.568	3660	3669	3678	rBV	776389	2115535	71.80%	8.440%

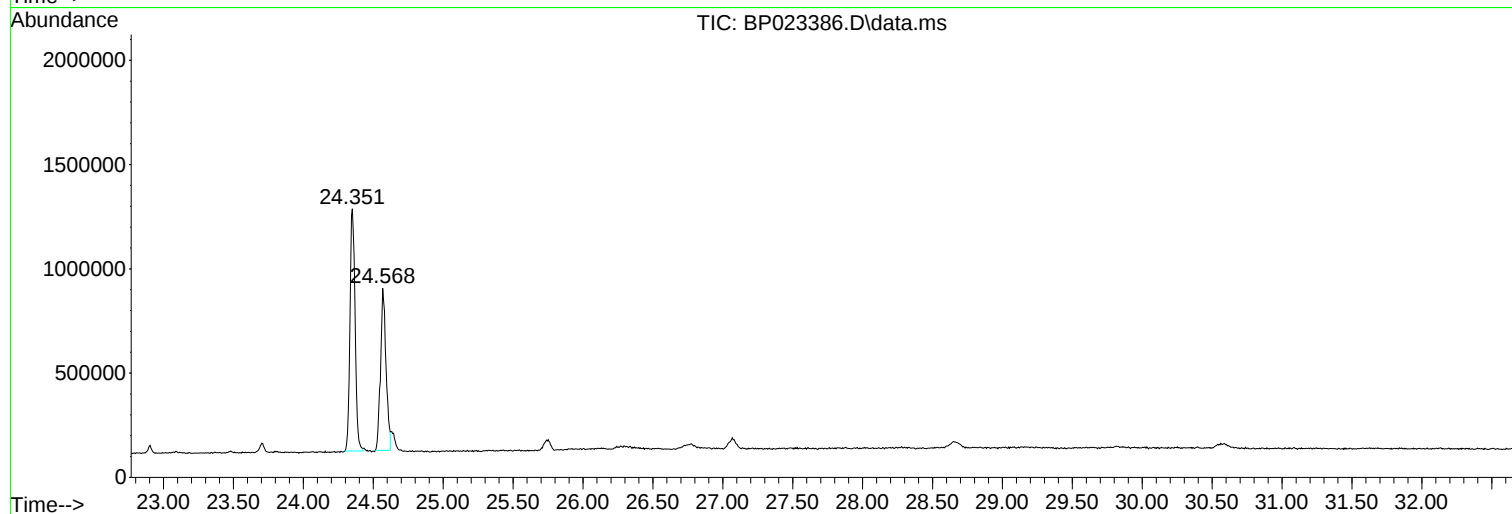
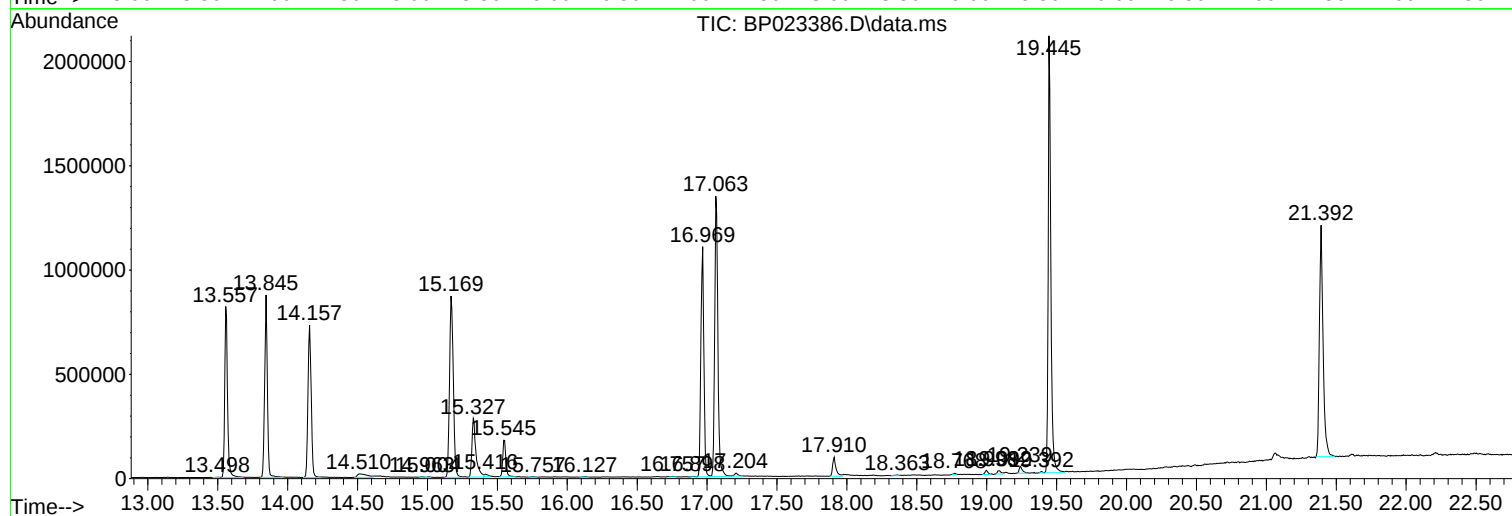
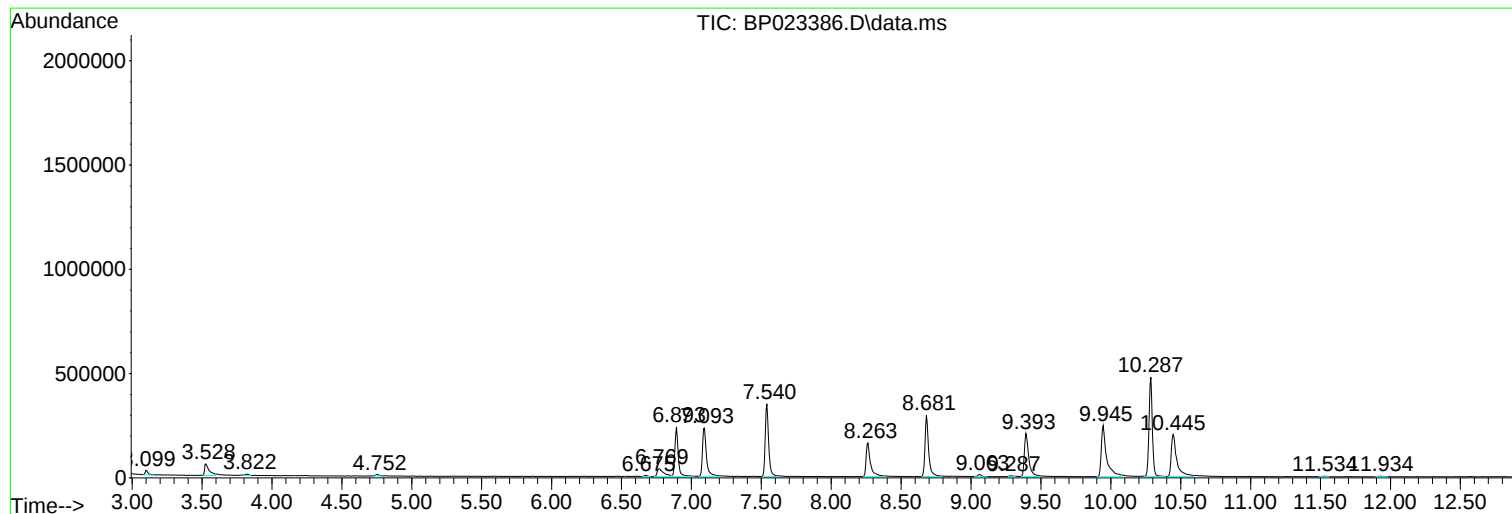
Sum of corrected areas: 25065743

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION

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



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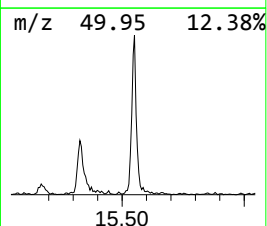
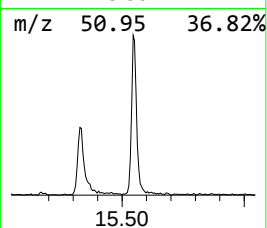
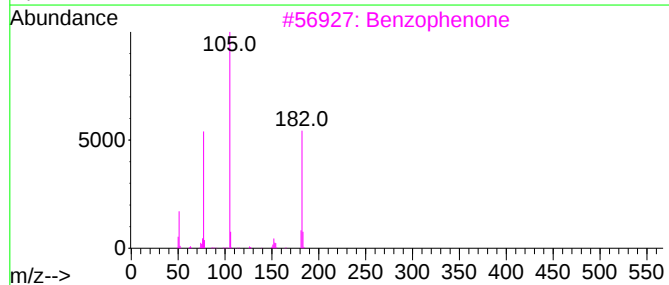
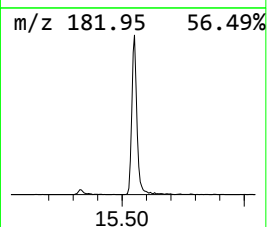
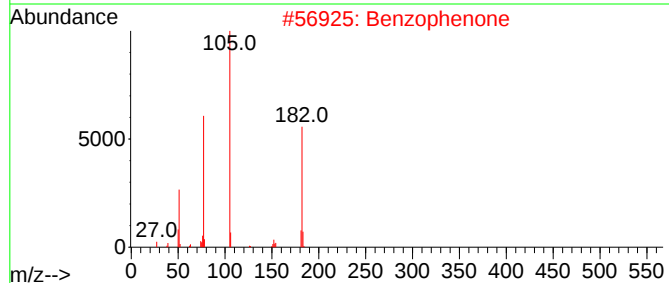
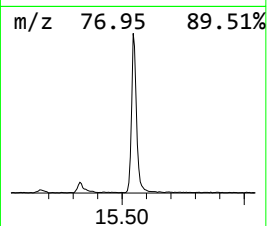
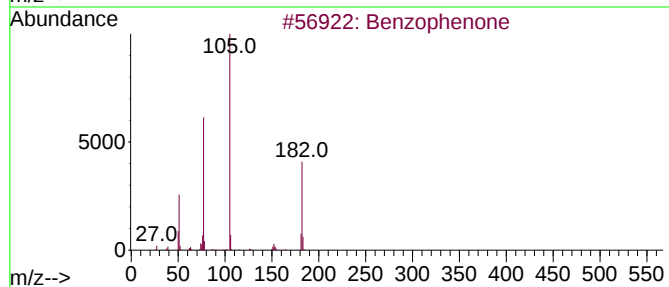
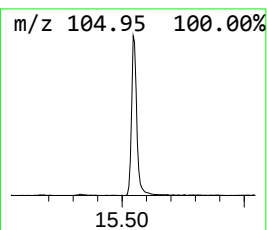
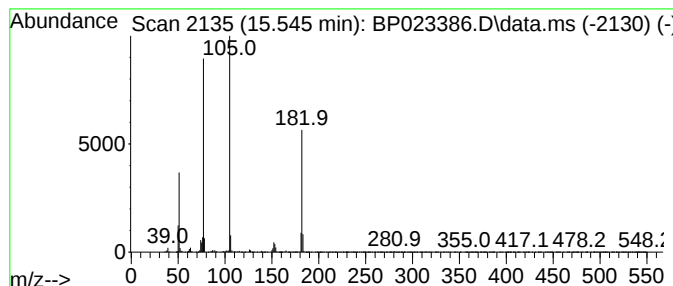
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION

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.
15.545	4.54 ng/ul	265691	Acenaphthene-d10	14.157

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	93
5			Benzophenone	182	C13H10O	000119-61-9	91



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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzophenone	15.545	4.5	ng/u1	265691	3	14.157	1171090	20.0