

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012025\
 Data File : BP023694.D
 Acq On : 21 Jan 2025 22:12
 Operator : RC/JU
 Sample : Q1108-19
 Misc :
 ALS Vial : 47 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 YE8H9

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-BP011425.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP023694.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.311	35	38	47	rVB	105078	139280	1.87%	0.161%
2	3.716	103	107	125	rBV	863078	1256673	16.89%	1.456%
3	4.052	160	164	169	rVB	35585	46587	0.63%	0.054%
4	4.958	313	318	326	rVB2	26972	45067	0.61%	0.052%
5	6.210	526	531	534	rVB5	6677	9373	0.13%	0.011%
6	6.963	651	659	674	rBV	1464453	2240681	30.12%	2.596%
7	7.134	681	688	701	rVB	1258263	1931066	25.96%	2.237%
8	7.340	715	723	732	rBV	1473225	2256528	30.33%	2.614%
9	7.446	736	741	744	rVB6	5042	9063	0.12%	0.010%
10	7.804	794	802	809	rBV	2211941	3386371	45.52%	3.923%
11	7.869	809	813	820	rVB4	19468	32204	0.43%	0.037%
12	8.487	911	918	932	rBV	1632258	2565890	34.49%	2.972%
13	8.581	932	934	941	rVB8	5622	12105	0.16%	0.014%
14	8.940	988	995	1006	rBV	1290043	2011192	27.03%	2.330%
15	9.116	1019	1025	1030	rBV7	9675	16158	0.22%	0.019%
16	9.398	1069	1073	1079	rVB2	25488	39564	0.53%	0.046%
17	9.504	1085	1091	1100	rBV	75221	111657	1.50%	0.129%
18	9.593	1103	1106	1107	rBV2	8431	9280	0.12%	0.011%
19	9.663	1110	1118	1126	rVV	737608	1176503	15.81%	1.363%
20	10.193	1199	1208	1230	rBV	1579883	2458781	33.05%	2.848%
21	10.369	1236	1238	1246	rVB7	8155	12527	0.17%	0.015%
22	10.575	1264	1273	1282	rBV	2516368	4369465	58.73%	5.062%
23	10.704	1287	1295	1306	rBV	1519407	2506916	33.70%	2.904%
24	11.110	1358	1364	1373	rVB2	32658	56462	0.76%	0.065%
25	12.010	1511	1517	1521	rBV9	8834	14236	0.19%	0.016%
26	12.163	1536	1543	1549	rBV	29164	49537	0.67%	0.057%
27	13.204	1716	1720	1725	rBV7	5112	11594	0.16%	0.013%
28	13.263	1725	1730	1734	rVV	19617	26012	0.35%	0.030%
29	13.387	1749	1751	1758	rVB8	7778	13802	0.19%	0.016%
30	13.828	1819	1826	1832	rBV	2178275	3091157	41.55%	3.581%
31	13.869	1832	1833	1842	rVB5	19071	27164	0.37%	0.031%
32	14.116	1865	1875	1884	rBV	2776933	4224208	56.78%	4.894%
33	14.345	1909	1914	1919	rVB8	7111	11579	0.16%	0.013%
34	14.422	1919	1927	1935	rBV	3181663	4874127	65.52%	5.646%
35	14.598	1950	1957	1970	rBV	765727	1234778	16.60%	1.430%
36	14.845	1997	1999	2005	rVB6	5969	10321	0.14%	0.012%

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Integration Parameters: LSCINT.P

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Peak Location: TOP

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP011425.MA.M
 Title : SVOA CALIBRATION

37	15.257	2064	2069	2073	rBV3	11873	19077	0.26%	0.022%
38	15.433	2091	2099	2111	rBV	3345177	4977092	66.90%	5.766%
39	15.539	2111	2117	2130	rBV	405184	587787	7.90%	0.681%
40	15.669	2136	2139	2145	rVB3	16034	26335	0.35%	0.031%
41	15.804	2156	2162	2173	rBV	974553	1373965	18.47%	1.592%
42	16.004	2191	2196	2202	rBV10	5374	9365	0.13%	0.011%
43	16.228	2232	2234	2240	rVB7	5709	9669	0.13%	0.011%
44	16.375	2254	2259	2268	rVB	57200	84542	1.14%	0.098%
45	17.004	2362	2366	2370	rVB3	16712	23077	0.31%	0.027%
46	17.169	2390	2394	2396	rBV5	8410	9427	0.13%	0.011%
47	17.227	2397	2404	2414	rVV	3590782	5237038	70.39%	6.067%
48	17.327	2415	2421	2433	rBV	4106848	5241719	70.46%	6.072%
49	17.639	2470	2474	2480	rVB7	23078	36308	0.49%	0.042%
50	17.933	2521	2524	2527	rBV5	5780	9299	0.12%	0.011%
51	18.098	2548	2552	2556	rBV5	18465	32556	0.44%	0.038%
52	18.169	2558	2564	2570	rBV	263281	408234	5.49%	0.473%
53	18.474	2613	2616	2621	rVB3	22134	30195	0.41%	0.035%
54	18.622	2637	2641	2646	rBV2	108917	133485	1.79%	0.155%
55	19.039	2707	2712	2721	rVB5	41919	82082	1.10%	0.095%
56	19.239	2742	2746	2752	rBV3	43242	60801	0.82%	0.070%
57	19.310	2755	2758	2762	rVV	31494	36419	0.49%	0.042%
58	19.480	2784	2787	2794	rBV4	59843	95548	1.28%	0.111%
59	19.680	2815	2821	2829	rBV	5224705	6254287	84.07%	7.245%
60	21.327	3098	3101	3109	rBV3	166831	281898	3.79%	0.327%
61	21.651	3150	3156	3167	rBV2	4124208	6474988	87.03%	7.501%
62	24.804	3684	3692	3713	rVB	2850569	7058814	94.88%	8.177%
63	25.021	3720	3729	3747	rVB2	2780944	7439567	100.00%	8.618%

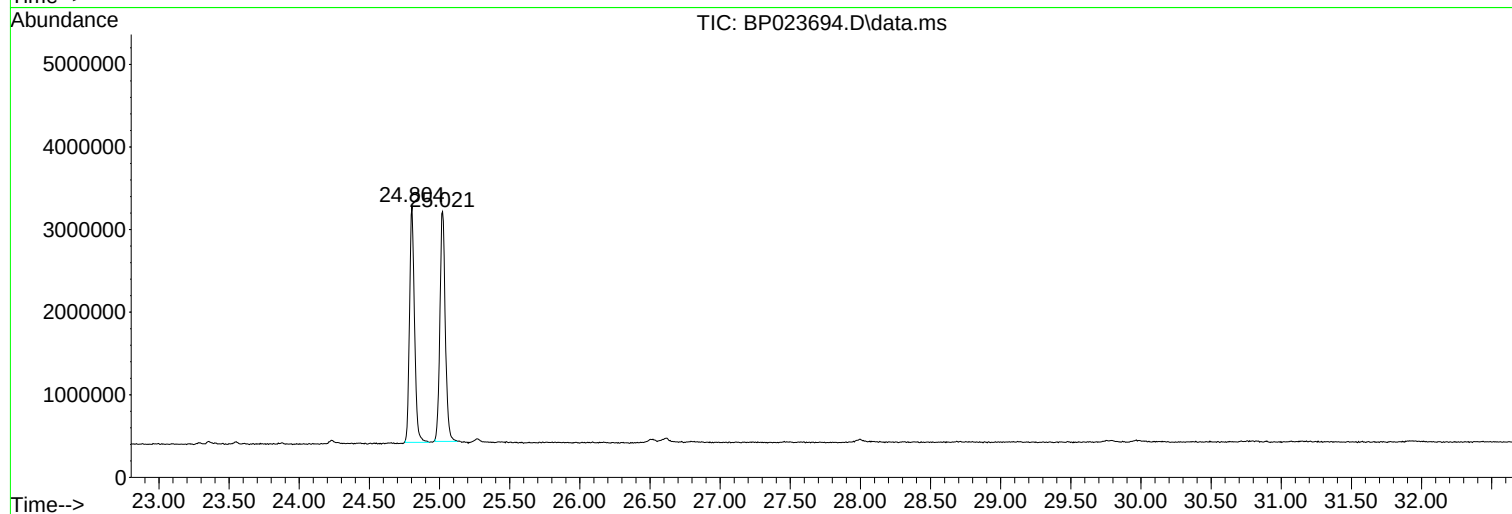
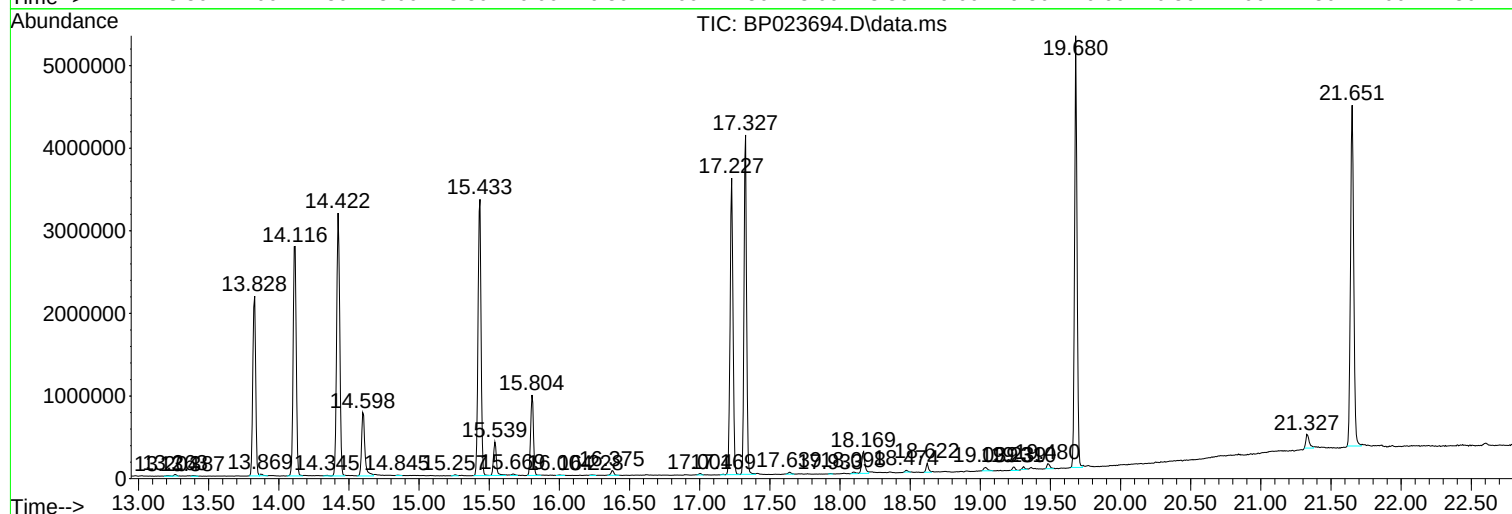
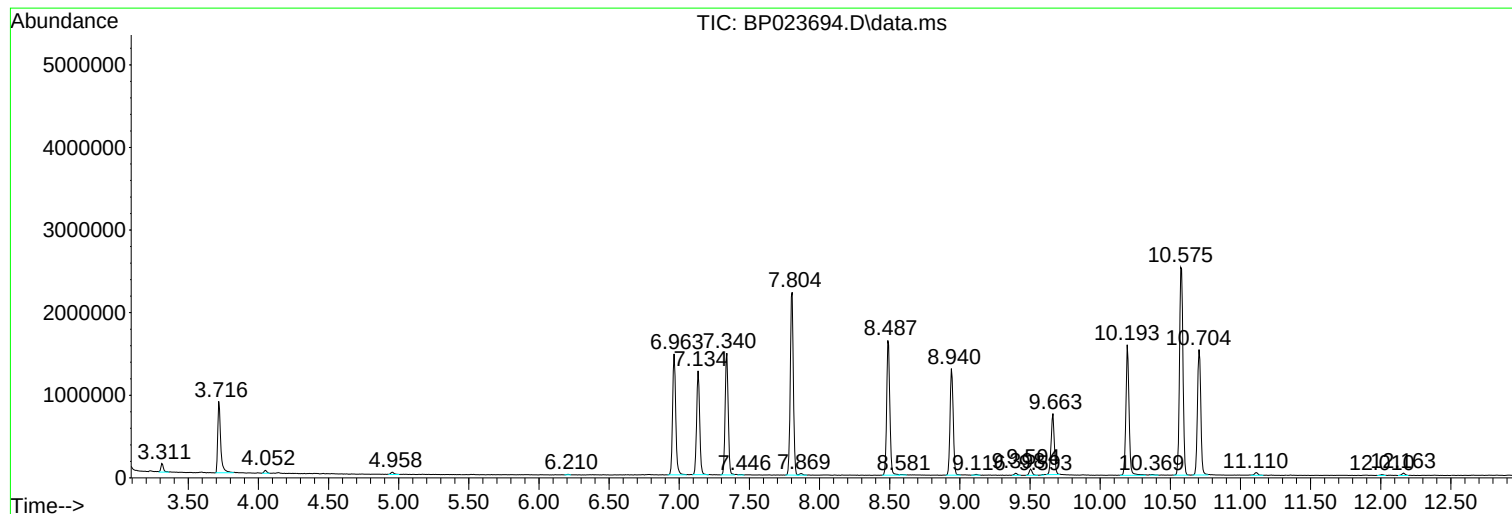
Sum of corrected areas: 86321482

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP011425.MA.M
Quant Title : SVOA CALIBRATION

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



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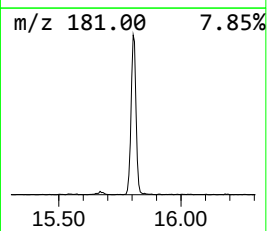
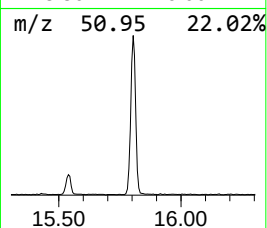
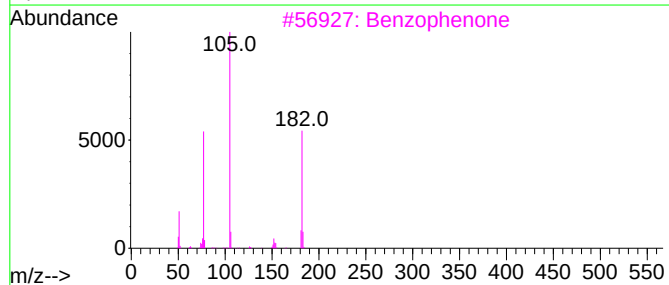
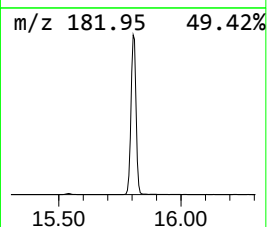
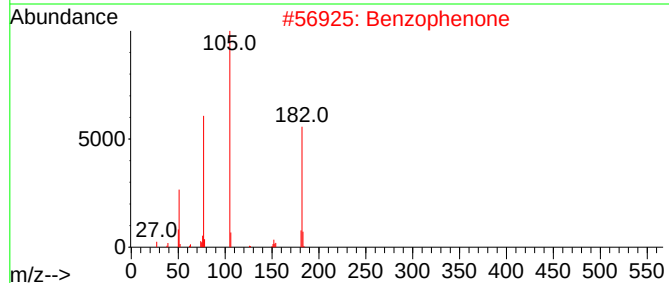
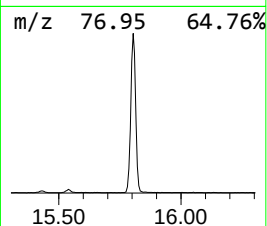
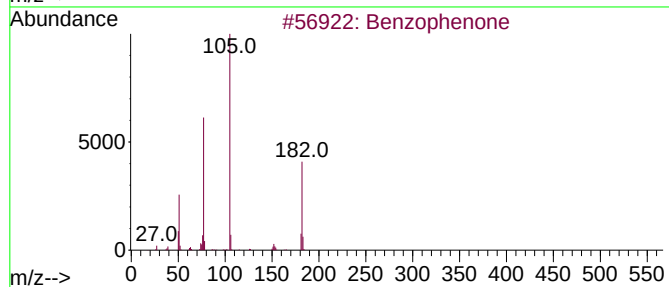
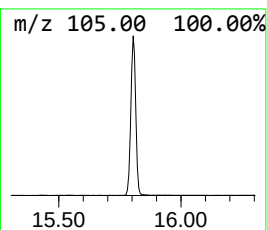
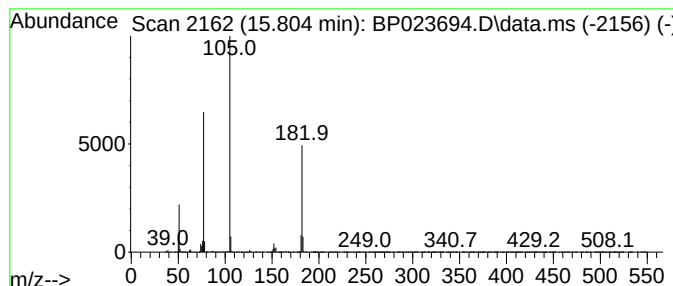
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP011425.MA.M
Quant Title : SVOA CALIBRATION

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.
15.804	5.64 ng/ul	1373970	Acenaphthene-d10	14.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	97
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 Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzophenone	15.804	5.6	ng/ul	1373970	3	14.422	4874130	20.0