

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101424\
 Data File : BP022392.D
 Acq On : 15 Oct 2024 16:58
 Operator : RC/JU
 Sample : P4349-09
 Misc :
 ALS Vial : 45 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EOAG8

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-BP100724.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP022392.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	27	rVB5	4072	5722	0.16%	0.019%
2	3.222	36	40	45	rVB	19430	24178	0.67%	0.080%
3	3.646	108	112	125	rBV	52367	90438	2.51%	0.301%
4	3.952	160	164	170	rVB4	6665	10308	0.29%	0.034%
5	4.111	190	191	198	rVB7	3394	4593	0.13%	0.015%
6	4.175	200	202	206	rVB4	3259	3659	0.10%	0.012%
7	4.322	223	227	234	rVB4	5337	8482	0.24%	0.028%
8	4.458	244	250	254	rBV4	5052	9336	0.26%	0.031%
9	4.505	254	258	264	rVB2	14376	22933	0.64%	0.076%
10	4.740	293	298	301	rBV6	2335	3948	0.11%	0.013%
11	4.864	313	319	325	rBV6	2842	6818	0.19%	0.023%
12	5.546	429	435	441	rBV9	3671	6970	0.19%	0.023%
13	5.734	459	467	476	rBV4	10033	19936	0.55%	0.066%
14	5.987	506	510	516	rVB9	2490	4972	0.14%	0.017%
15	6.105	520	530	538	rVB2	29191	63622	1.76%	0.212%
16	6.528	598	602	605	rVV5	2950	4158	0.12%	0.014%
17	6.569	605	609	615	rVB3	7890	11108	0.31%	0.037%
18	6.840	649	655	656	rBV6	2837	4138	0.11%	0.014%
19	6.887	656	663	676	rVB	179663	299280	8.29%	0.995%
20	7.034	682	688	698	rBV	275049	429425	11.90%	1.428%
21	7.240	715	723	734	rBV	395703	624951	17.32%	2.078%
22	7.693	793	800	810	rBV	400936	652909	18.09%	2.171%
23	7.805	814	819	824	rVB8	2939	5302	0.15%	0.018%
24	8.393	912	919	937	rBV	392845	652101	18.07%	2.168%
25	8.834	986	994	1012	rBV	358497	603427	16.72%	2.007%
26	9.246	1061	1064	1071	rVB6	4814	8284	0.23%	0.028%
27	9.393	1082	1089	1095	rBV6	2836	5927	0.16%	0.020%
28	9.546	1107	1115	1122	rBV	312294	534272	14.81%	1.777%
29	10.087	1199	1207	1220	rBV	491833	867752	24.05%	2.885%
30	10.451	1261	1269	1280	rVB	527070	859237	23.81%	2.857%
31	10.587	1285	1292	1312	rBV	270231	486365	13.48%	1.617%
32	11.269	1401	1408	1414	rBV9	3677	8333	0.23%	0.028%
33	12.040	1531	1539	1546	rBV2	7663	16439	0.46%	0.055%
34	13.716	1817	1824	1833	rBV	936984	1321151	36.61%	4.393%
35	13.998	1863	1872	1882	rBV	912005	1433652	39.73%	4.767%
36	14.304	1915	1924	1932	rBV	762671	1287450	35.68%	4.281%

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Peak Location: TOP

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Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
 Title : SVOA CALIBRATION

37	14.522	1956	1961	1979	rBV	163727	309409	8.57%	1.029%
38	14.681	1984	1988	1995	rVB8	4933	8928	0.25%	0.030%
39	14.739	1995	1998	2000	rBV3	3528	5272	0.15%	0.018%
40	15.310	2088	2095	2103	rBV	1424023	2032498	56.32%	6.758%
41	15.457	2112	2120	2128	rBV	587654	712379	19.74%	2.369%
42	15.563	2134	2138	2142	rVB3	7723	10756	0.30%	0.036%
43	15.692	2152	2160	2169	rBV	183692	326528	9.05%	1.086%
44	15.892	2187	2194	2195	rBV7	2399	4455	0.12%	0.015%
45	16.034	2213	2218	2225	rBV	45176	73397	2.03%	0.244%
46	16.128	2230	2234	2239	rVB8	2936	4961	0.14%	0.016%
47	16.263	2251	2257	2263	rBV	50178	74534	2.07%	0.248%
48	16.504	2294	2298	2302	rBV7	6715	11012	0.31%	0.037%
49	16.781	2340	2345	2349	rBV7	3779	7508	0.21%	0.025%
50	16.892	2362	2364	2369	rVB5	4090	5413	0.15%	0.018%
51	16.975	2375	2378	2384	rVB8	4456	5696	0.16%	0.019%
52	17.116	2394	2402	2409	rBV	1179764	1747372	48.42%	5.810%
53	17.210	2412	2418	2429	rVV	1716449	2447324	67.82%	8.138%
54	17.498	2464	2467	2468	rBV3	4318	4609	0.13%	0.015%
55	17.904	2533	2536	2543	rVB8	10165	16304	0.45%	0.054%
56	18.045	2555	2560	2570	rBV	66434	118183	3.27%	0.393%
57	18.551	2644	2646	2651	rVB5	8786	10392	0.29%	0.035%
58	18.933	2706	2711	2716	rVB2	19859	30273	0.84%	0.101%
59	19.127	2741	2744	2752	rVB4	24684	34394	0.95%	0.114%
60	19.216	2754	2759	2763	rBV	26893	52171	1.45%	0.173%
61	19.375	2782	2786	2795	rBV6	34711	59741	1.66%	0.199%
62	19.533	2811	2813	2816	rVB3	12267	11576	0.32%	0.038%
63	19.592	2817	2823	2831	rBV	2439782	3294300	91.29%	10.954%
64	21.539	3148	3154	3163	rBV	1515328	2257719	62.56%	7.507%
65	24.604	3666	3675	3687	rVB	1347438	3608656	100.00%	11.999%
66	24.845	3707	3716	3728	rVB	1003685	2386253	66.13%	7.935%

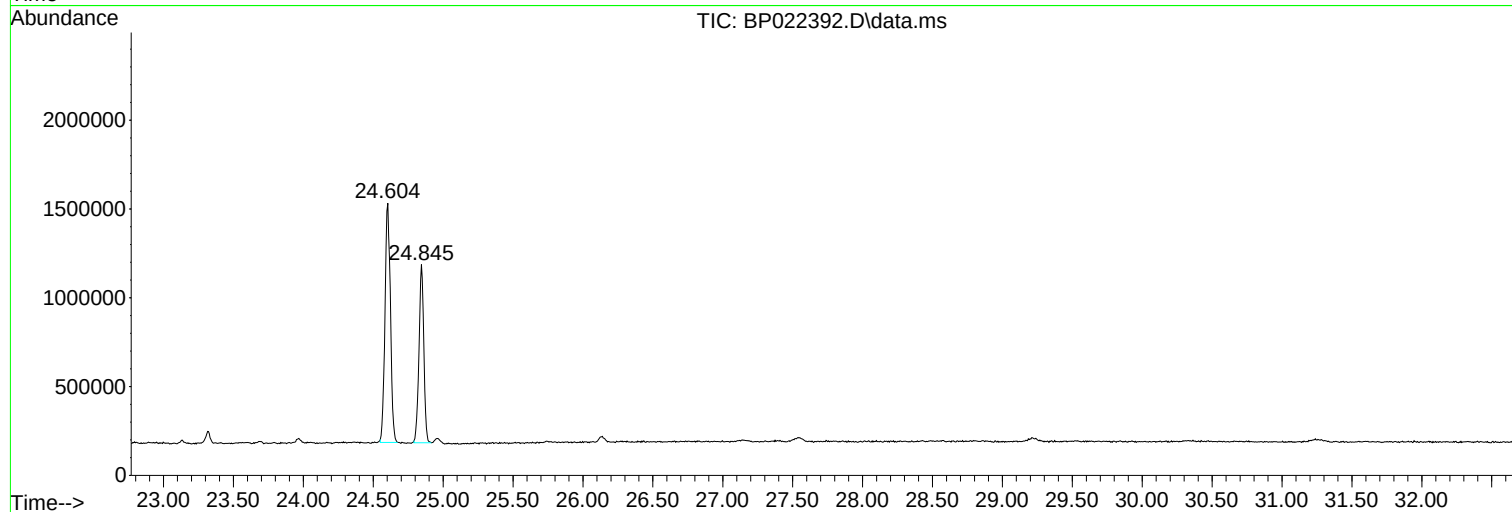
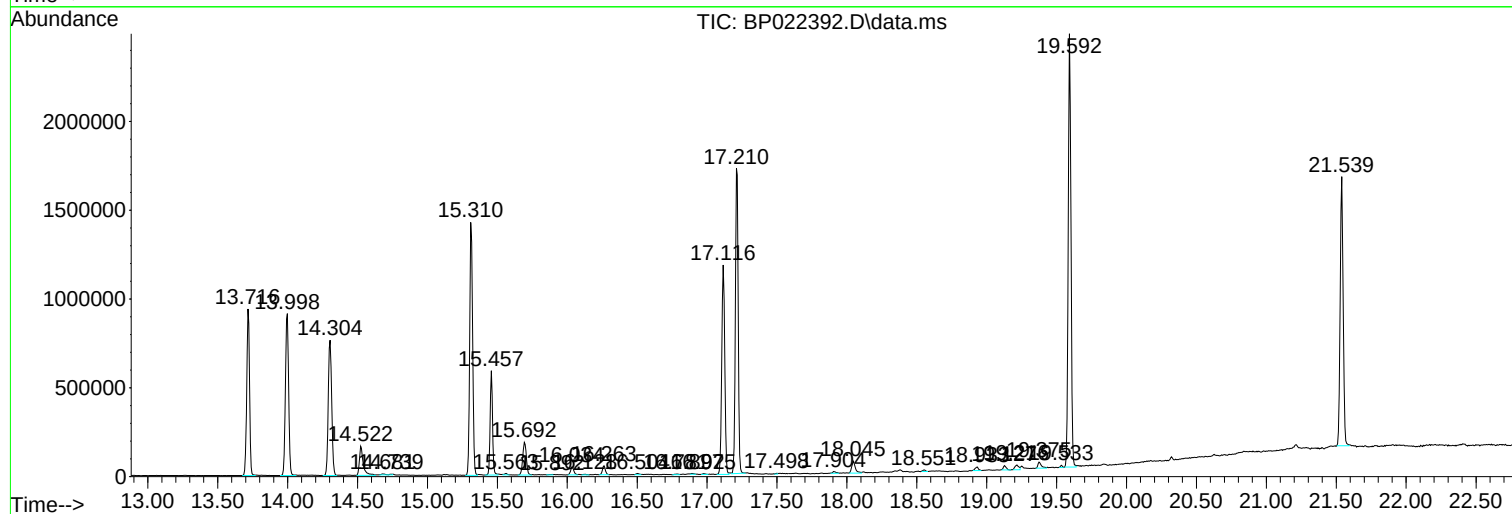
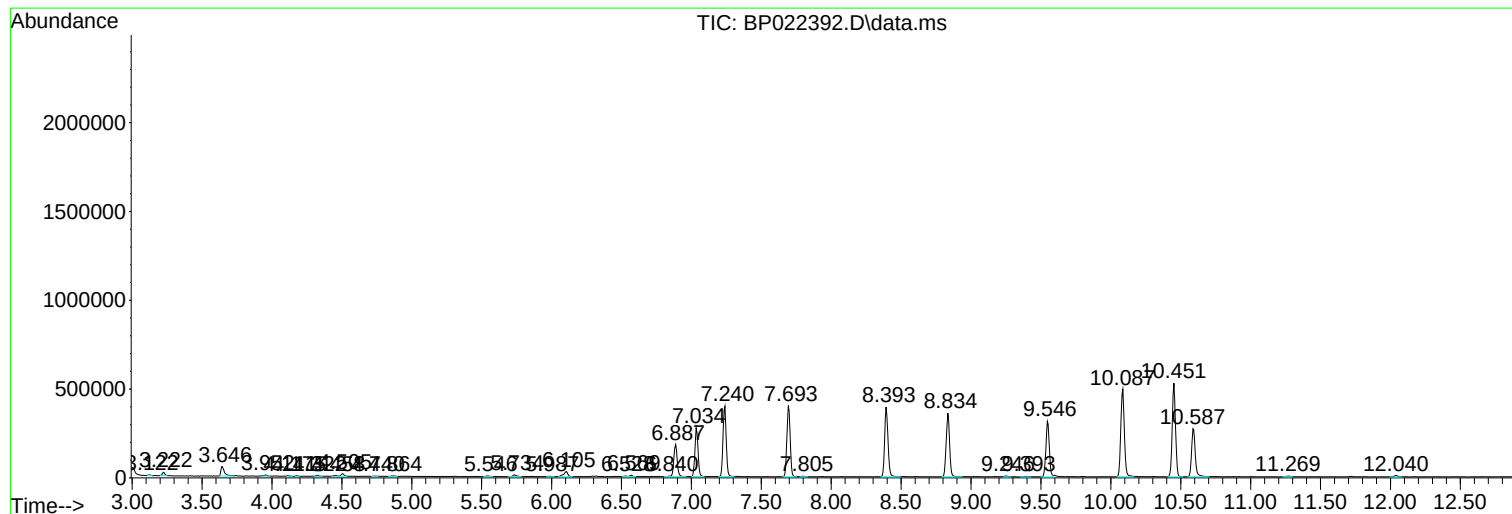
Sum of corrected areas: 30073589

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION

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



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Instrument :
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ClientSampleId :
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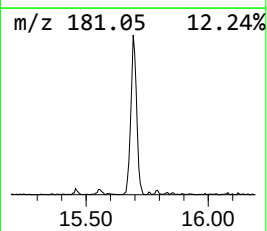
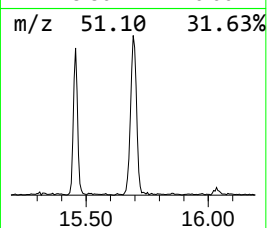
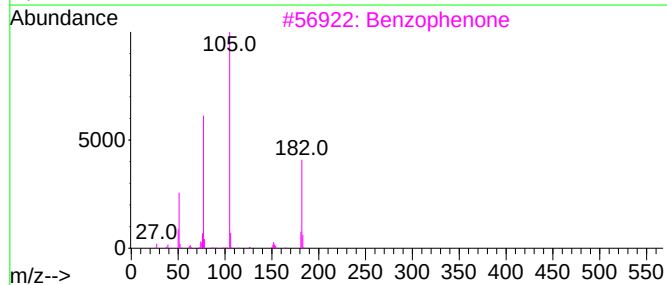
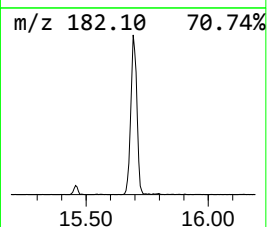
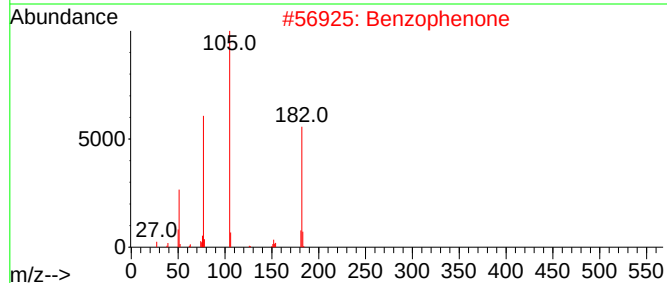
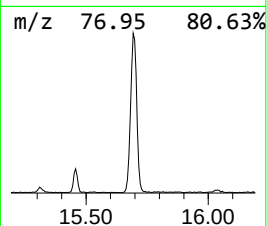
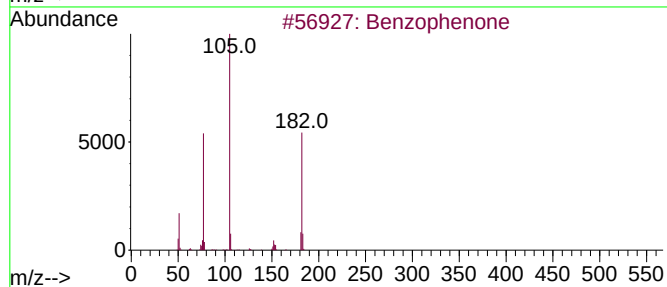
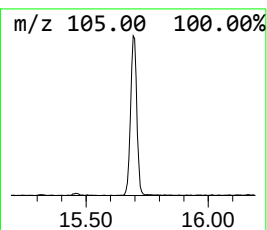
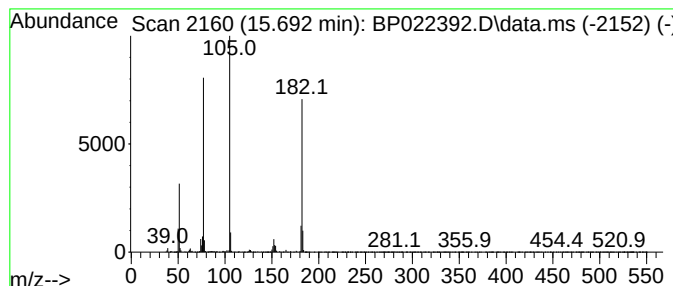
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION

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.
15.692	5.07 ng/ul	326528	Acenaphthene-d10	14.304

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	94
4			Benzophenone	182	C13H10O	000119-61-9	91
5			1,2,4-Trioxolane, 3,3,5-triphenyl-	304	C20H16O3	023246-12-0	78



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TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzophenone	15.692	5.1	ng/u1	326528	3	14.304	1287450	20.0