

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101424\
 Data File : BP022398.D
 Acq On : 15 Oct 2024 21:02
 Operator : RC/JU
 Sample : P4349-17
 Misc :
 ALS Vial : 51 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EOAG1

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: BP022398.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.223	37	40	47	rVB	15656	19960	0.57%	0.075%
2	3.646	108	112	124	rBV	43604	73779	2.11%	0.277%
3	3.952	161	164	170	rBV3	4246	6163	0.18%	0.023%
4	4.505	252	258	266	rVB7	4489	7782	0.22%	0.029%
5	6.887	656	663	676	rVB	91685	163008	4.67%	0.612%
6	7.040	681	689	699	rBV	217431	363273	10.41%	1.364%
7	7.240	715	723	735	rBV	293852	490342	14.05%	1.841%
8	7.693	792	800	809	rBV	304987	491382	14.08%	1.845%
9	8.399	912	920	935	rBV	260248	442576	12.69%	1.662%
10	8.834	985	994	1007	rBV	307331	531073	15.22%	1.994%
11	9.246	1058	1064	1069	rBV3	4953	7657	0.22%	0.029%
12	9.393	1084	1089	1095	rBB3	7941	12682	0.36%	0.048%
13	9.546	1104	1115	1125	rBV	270519	479195	13.73%	1.799%
14	10.081	1198	1206	1223	rBV	443646	769998	22.07%	2.891%
15	10.452	1260	1269	1277	rBV	379573	684215	19.61%	2.569%
16	10.593	1283	1293	1307	rBV	206744	388028	11.12%	1.457%
17	11.375	1415	1426	1430	rBV	2241	6684	0.19%	0.025%
18	12.034	1531	1538	1548	rBV3	6807	14291	0.41%	0.054%
19	13.704	1816	1822	1835	rBV	953456	1247918	35.77%	4.685%
20	13.810	1838	1840	1848	rVB7	2008	4279	0.12%	0.016%
21	13.998	1862	1872	1882	rBV	856483	1361312	39.02%	5.111%
22	14.310	1916	1925	1934	rBV	648587	1068933	30.64%	4.013%
23	14.540	1956	1964	1978	rBV2	80243	199785	5.73%	0.750%
24	14.692	1985	1990	1994	rBV7	4364	7023	0.20%	0.026%
25	14.745	1995	1999	2005	rVB8	5243	8018	0.23%	0.030%
26	15.316	2089	2096	2104	rBV	1249622	1934291	55.44%	7.262%
27	15.369	2104	2105	2111	rVV6	3769	5927	0.17%	0.022%
28	15.445	2112	2118	2130	rVV	580871	709665	20.34%	2.664%
29	15.557	2133	2137	2142	rVB5	5646	9858	0.28%	0.037%
30	15.698	2155	2161	2166	rBV	126199	179317	5.14%	0.673%
31	16.039	2212	2219	2229	rBV	29018	50491	1.45%	0.190%
32	16.122	2229	2233	2237	rVB7	2343	4533	0.13%	0.017%
33	16.716	2328	2334	2337	rBV7	3474	6267	0.18%	0.024%
34	17.110	2395	2401	2407	rBV	1129700	1526379	43.75%	5.731%
35	17.157	2407	2409	2411	rVV3	7848	8851	0.25%	0.033%
36	17.204	2411	2417	2428	rVB2	1604638	2449047	70.19%	9.195%

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37	17.498	2462	2467	2468	rBV5	5145	7587	0.22%	0.028%
38	18.045	2555	2560	2565	rBV	51480	74173	2.13%	0.278%
39	18.539	2642	2644	2649	rBV5	5906	8304	0.24%	0.031%
40	18.916	2704	2708	2712	rBV7	7850	15823	0.45%	0.059%
41	19.122	2739	2743	2749	rVB3	21940	32555	0.93%	0.122%
42	19.204	2753	2757	2761	rBV	23074	37788	1.08%	0.142%
43	19.369	2782	2785	2790	rBV6	24948	34869	1.00%	0.131%
44	19.522	2808	2811	2814	rBV4	9711	10318	0.30%	0.039%
45	19.586	2815	2822	2832	rBV	2194410	3191279	91.47%	11.982%
46	21.533	3147	3153	3166	rBV2	1206046	1929253	55.30%	7.243%
47	24.610	3665	3676	3689	rVB	1218564	3488971	100.00%	13.099%
48	24.827	3704	3713	3727	rVB	734342	2069797	59.32%	7.771%

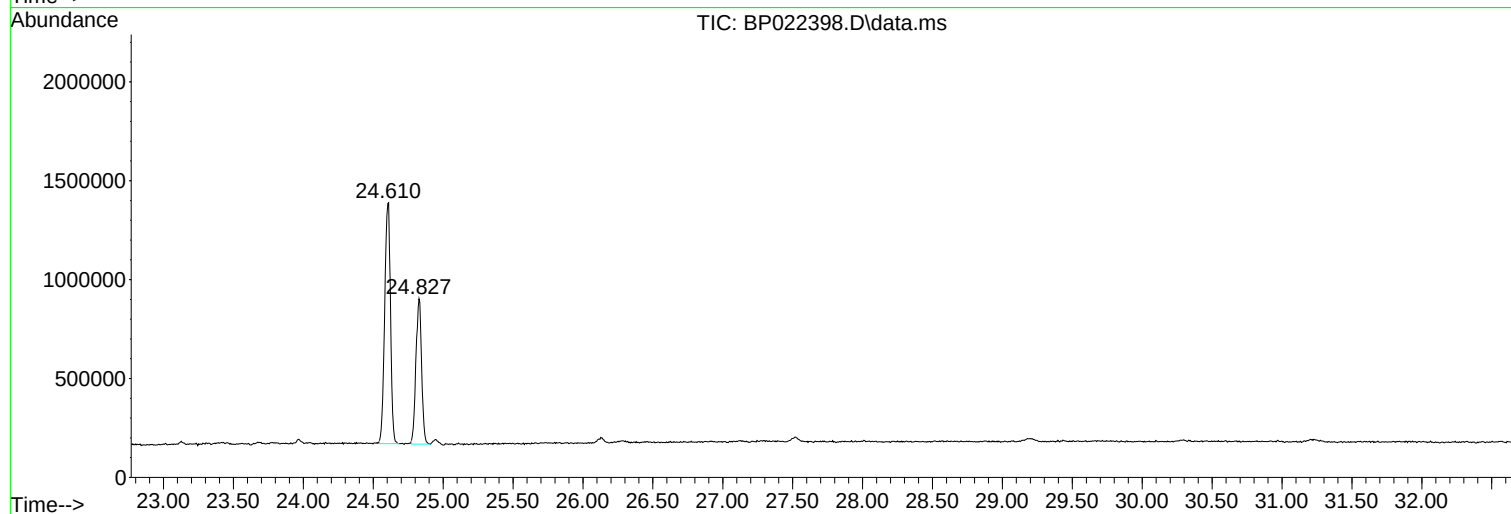
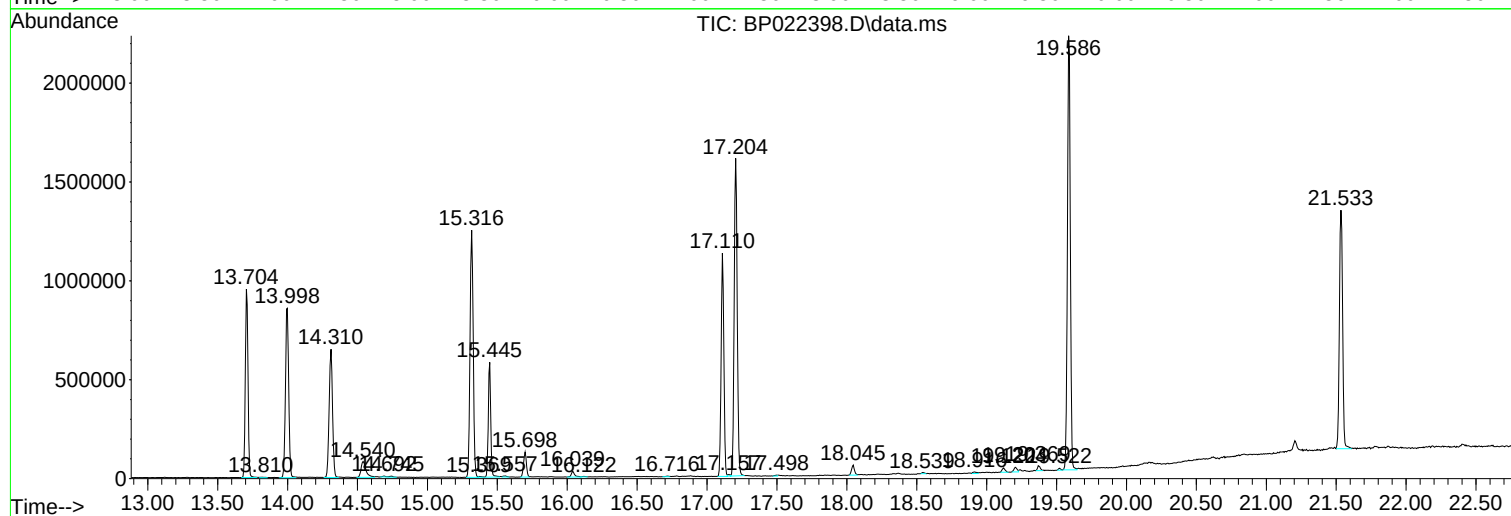
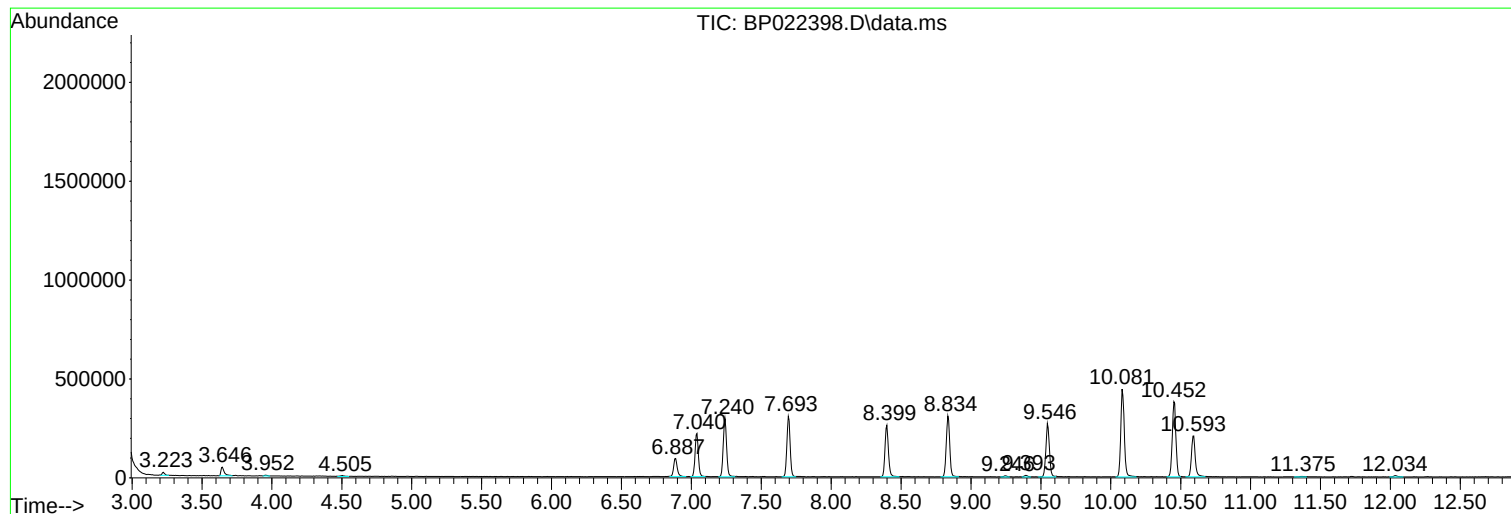
Sum of corrected areas: 26634699

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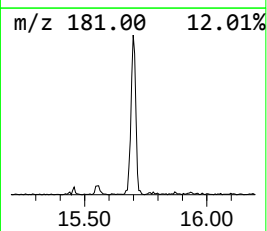
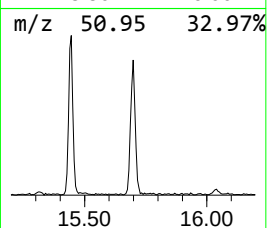
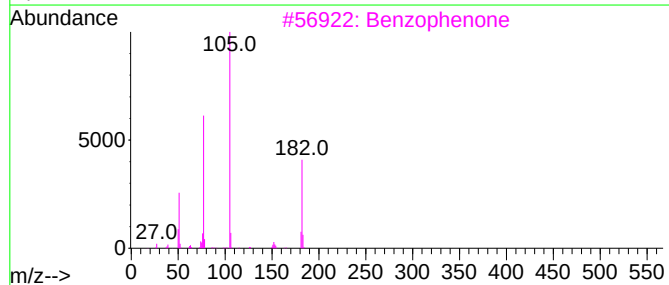
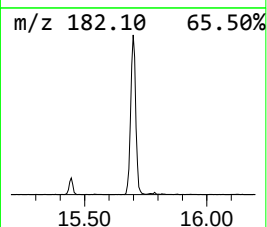
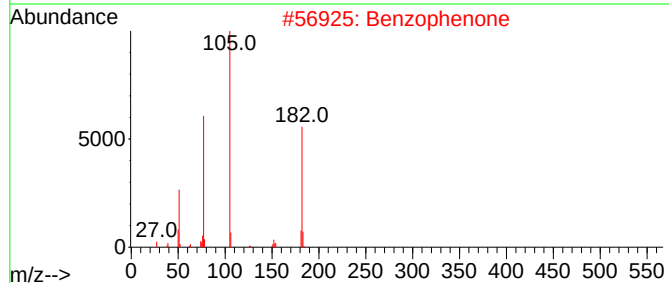
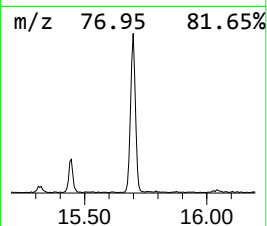
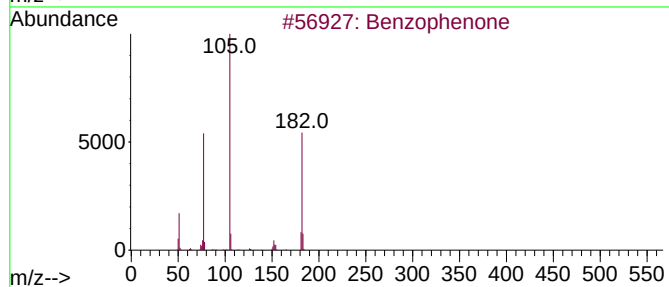
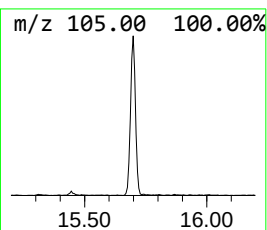
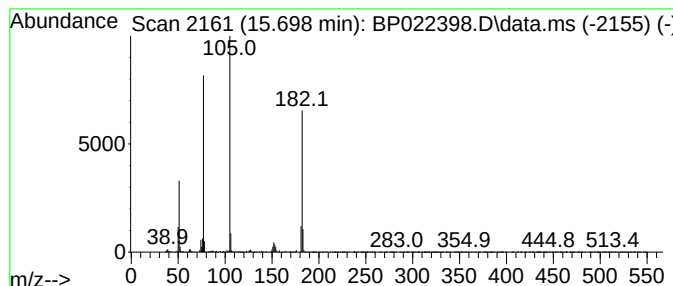
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION

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.
15.698	3.36 ng/ul	179317	Acenaphthene-d10	14.310

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	91
4			Benzophenone	182	C13H10O	000119-61-9	91
5			Benzophenone	182	C13H10O	000119-61-9	80



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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\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzophenone	15.698	3.4	ng/u1	179317	3	14.310	1068930	20.0