

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111324\
 Data File : BP023101.D
 Acq On : 14 Nov 2024 10:19
 Operator : RC/JU
 Sample : P4724-10
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BH9F2

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-BP110924.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP023101.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.140	22	26	30	rBV3	9149	12232	0.63%	0.069%
2	3.599	99	104	107	rBV7	3859	6894	0.36%	0.039%
3	3.858	145	148	154	rBV2	3455	5632	0.29%	0.032%
4	6.787	640	646	663	rBV	130045	269759	13.95%	1.531%
5	6.928	663	670	683	rVB	122630	200639	10.37%	1.138%
6	7.128	697	704	720	rBV	160878	272503	14.09%	1.546%
7	7.581	774	781	792	rBV	279702	477333	24.68%	2.708%
8	8.293	896	902	920	rBV	162602	327561	16.93%	1.859%
9	8.722	968	975	993	rBV	156540	298034	15.41%	1.691%
10	9.111	1034	1041	1047	rVB9	2746	5844	0.30%	0.033%
11	9.305	1070	1074	1081	rBV8	2441	5092	0.26%	0.029%
12	9.434	1089	1096	1112	rBV	137314	246547	12.75%	1.399%
13	9.981	1181	1189	1205	rBV	161168	377974	19.54%	2.145%
14	10.322	1236	1247	1257	rBV	364515	648402	33.52%	3.679%
15	10.481	1266	1274	1294	rBV	142283	331028	17.11%	1.878%
16	11.746	1486	1489	1494	rVB6	2112	3486	0.18%	0.020%
17	11.940	1517	1522	1526	rBV6	3087	6042	0.31%	0.034%
18	12.799	1663	1668	1671	rBV7	1336	2171	0.11%	0.012%
19	12.999	1697	1702	1707	rBV5	4040	5994	0.31%	0.034%
20	13.610	1797	1806	1819	rBV	490544	612886	31.68%	3.478%
21	13.887	1844	1853	1861	rBV	518916	765139	39.55%	4.342%
22	14.093	1883	1888	1892	rVB5	5018	6349	0.33%	0.036%
23	14.193	1898	1905	1914	rBV	648343	1029300	53.21%	5.840%
24	14.499	1949	1957	1965	rBV2	38912	119987	6.20%	0.681%
25	14.663	1983	1985	1991	rVB6	3933	5637	0.29%	0.032%
26	15.010	2039	2044	2046	rBV4	2475	3103	0.16%	0.018%
27	15.040	2046	2049	2050	rVV4	2463	2660	0.14%	0.015%
28	15.057	2050	2052	2056	rVB5	3800	5094	0.26%	0.029%
29	15.204	2069	2077	2088	rBV	631296	1056991	54.64%	5.998%
30	15.369	2100	2105	2116	rBV	58675	108119	5.59%	0.613%
31	15.451	2116	2119	2123	rVB6	3876	6176	0.32%	0.035%
32	15.581	2134	2141	2151	rBV	217883	344626	17.82%	1.955%
33	15.787	2173	2176	2181	rBV7	1611	3420	0.18%	0.019%
34	15.887	2192	2193	2196	rBV3	1688	1960	0.10%	0.011%
35	15.940	2199	2202	2206	rVB4	3925	5181	0.27%	0.029%
36	16.122	2229	2233	2236	rBV6	3736	5512	0.28%	0.031%

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Peak Location: TOP

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110924.MA.M
 Title : SVOA CALIBRATION

37	16.157	2236	2239	2243	rVB3	4431	5851	0.30%	0.033%
38	16.545	2299	2305	2306	rBV6	2414	4021	0.21%	0.023%
39	16.763	2337	2342	2346	rBV4	3775	6413	0.33%	0.036%
40	16.934	2366	2371	2375	rBV7	1877	2769	0.14%	0.016%
41	16.992	2375	2381	2392	rVV	1007062	1394702	72.10%	7.914%
42	17.104	2392	2400	2415	rVV	693379	1253249	64.79%	7.111%
43	17.310	2430	2435	2441	rVB9	3024	6862	0.35%	0.039%
44	17.516	2467	2470	2475	rVB6	4092	5354	0.28%	0.030%
45	17.939	2537	2542	2548	rBV4	31381	56000	2.89%	0.318%
46	18.234	2587	2592	2598	rBV6	12795	22563	1.17%	0.128%
47	18.387	2615	2618	2622	rBV5	3656	4394	0.23%	0.025%
48	18.675	2663	2667	2669	rBV4	2913	4342	0.22%	0.025%
49	18.810	2684	2690	2697	rVB5	8664	17927	0.93%	0.102%
50	18.898	2703	2705	2710	rVB6	5478	6685	0.35%	0.038%
51	19.039	2725	2729	2732	rBV4	12805	17409	0.90%	0.099%
52	19.110	2737	2741	2744	rVV3	9670	14554	0.75%	0.083%
53	19.275	2765	2769	2772	rBV5	9303	13158	0.68%	0.075%
54	19.486	2799	2805	2814	rBV	1217559	1601904	82.81%	9.090%
55	21.104	3076	3080	3098	rBV5	60481	196921	10.18%	1.117%
56	21.433	3130	3136	3149	rBV	1199778	1769212	91.46%	10.039%
57	24.416	3635	3643	3659	rVB	693076	1699544	87.86%	9.644%
58	24.657	3673	3684	3700	rVB	607107	1934439	100.00%	10.976%

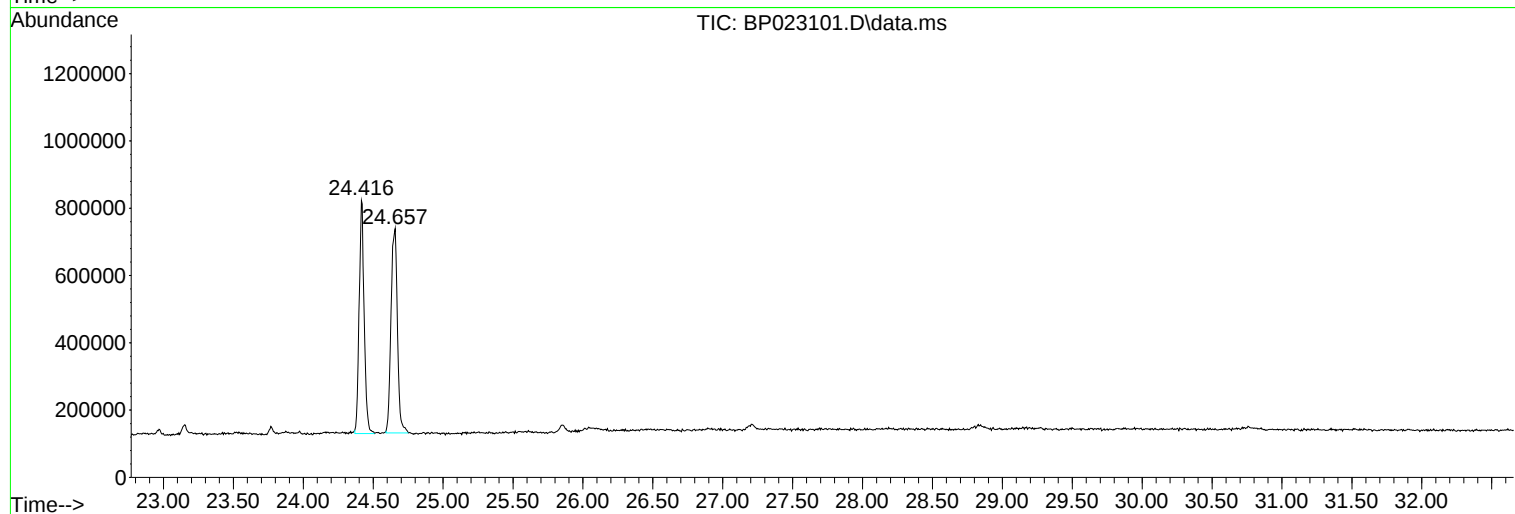
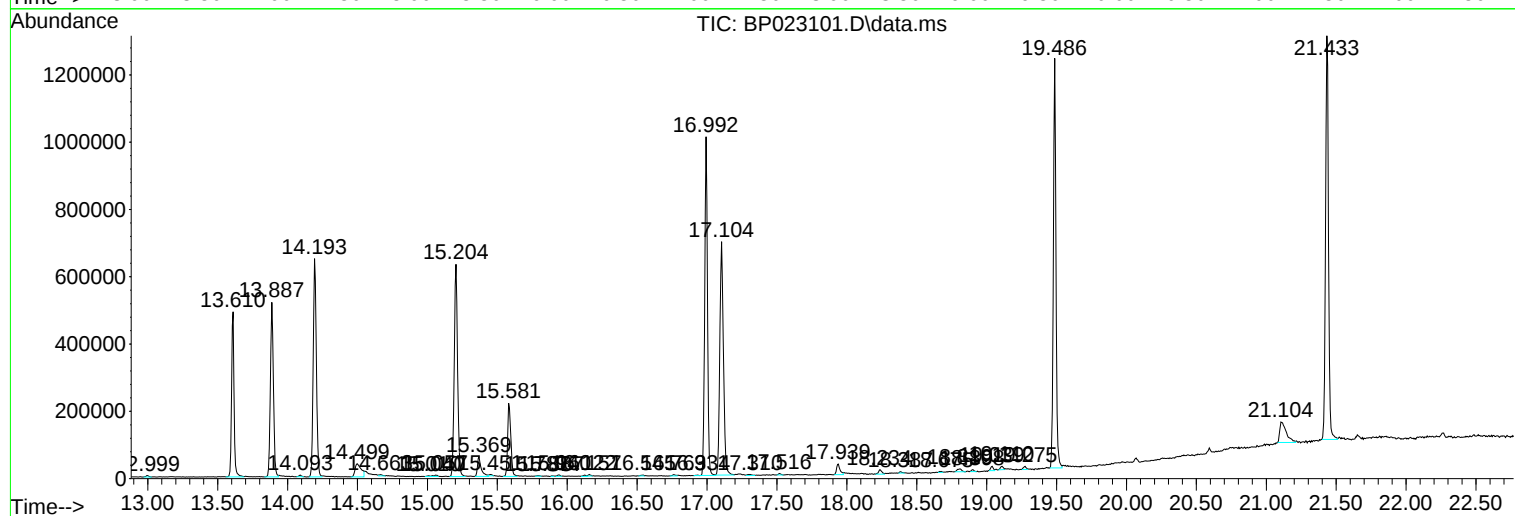
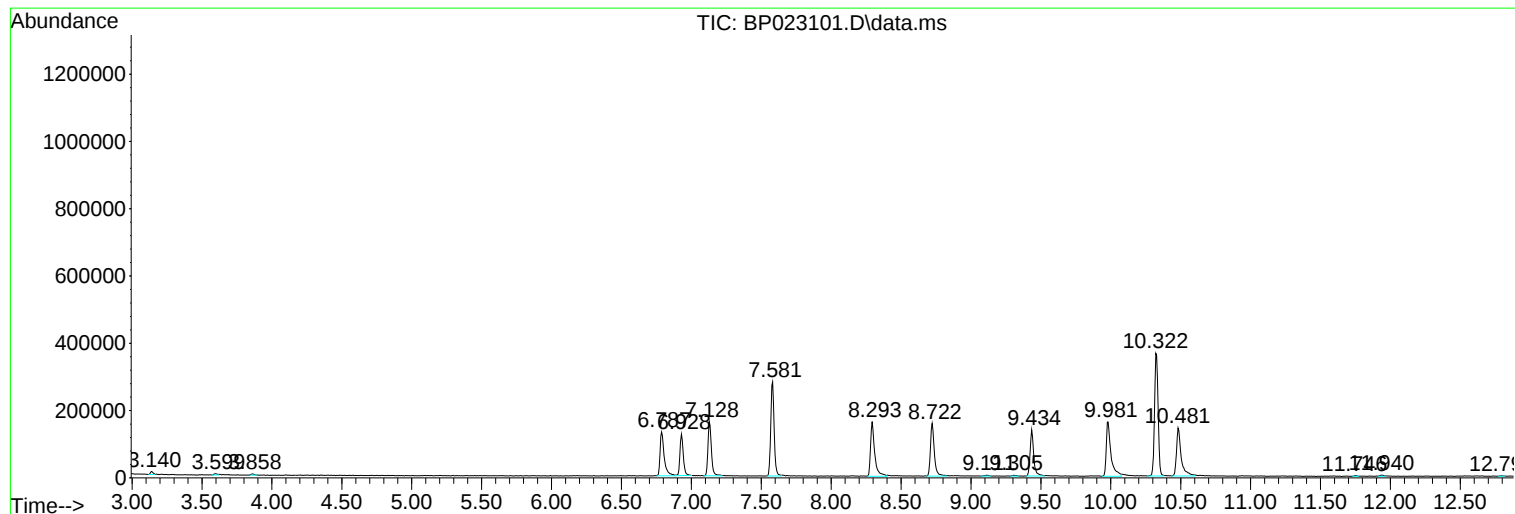
Sum of corrected areas: 17623580

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110924.MA.M
Quant Title : SVOA CALIBRATION

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



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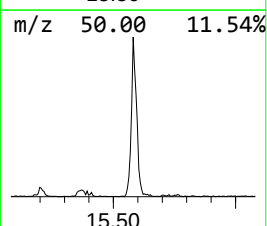
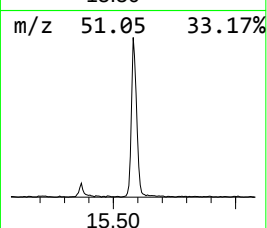
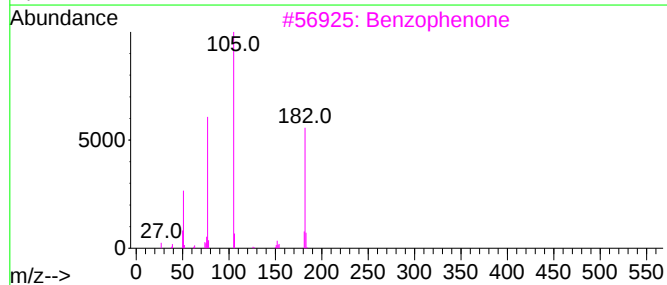
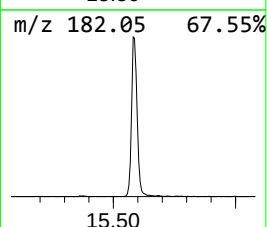
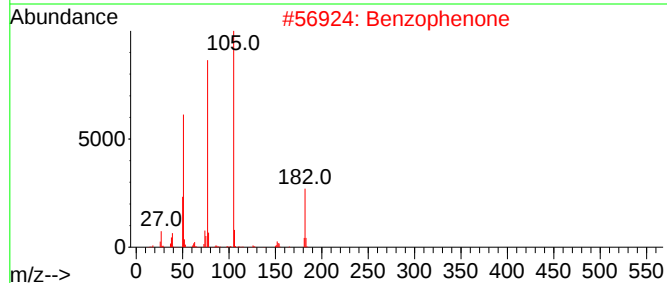
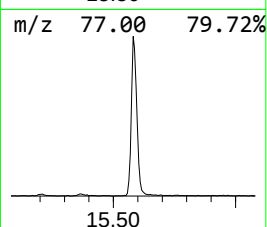
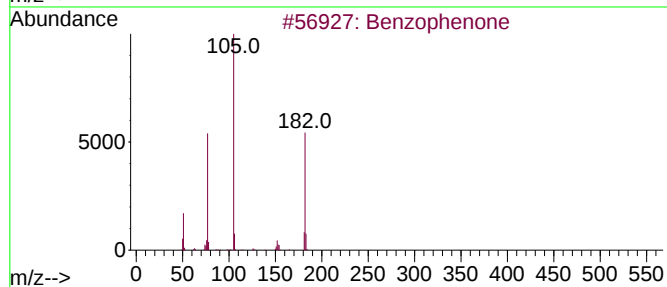
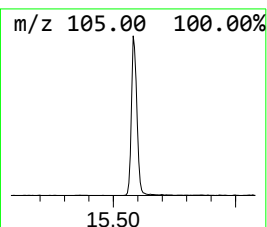
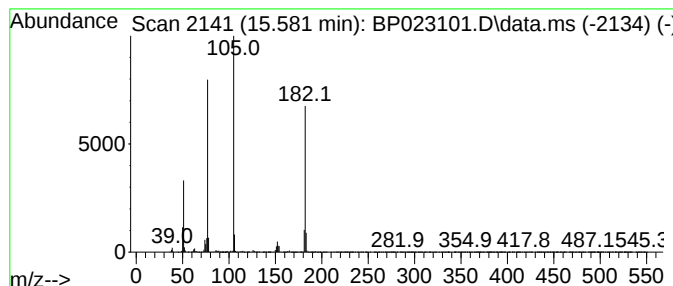
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110924.MA.M
Quant Title : SVOA CALIBRATION

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.
15.581	6.70 ng/ul	344626	Acenaphthene-d10	14.193

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	97
2			Benzophenone	182	C13H10O	000119-61-9	97
3			Benzophenone	182	C13H10O	000119-61-9	96
4			Benzophenone	182	C13H10O	000119-61-9	94
5			Benzophenone	182	C13H10O	000119-61-9	91



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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	15.581	6.7	ng/u1	344626	3	14.193	1029300	20.0