

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121024\
 Data File : BP023364.D
 Acq On : 11 Dec 2024 03:21
 Operator : RC/JU
 Sample : P5111-11
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BG460

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: BP023364.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.105	16	20	34	rVB	24141	39310	1.26%	0.149%
2	3.534	90	93	111	rBV	29864	71366	2.28%	0.271%
3	3.822	138	142	149	rVB6	4999	7873	0.25%	0.030%
4	4.246	210	214	220	rBV	12389	18559	0.59%	0.070%
5	4.746	295	299	307	rVB2	9185	15187	0.48%	0.058%
6	6.675	621	627	633	rBV2	7845	12499	0.40%	0.047%
7	6.758	635	641	658	rBV	77313	195086	6.23%	0.741%
8	6.893	658	664	675	rVB	307258	483258	15.43%	1.835%
9	7.087	691	697	714	rBV	333997	614219	19.61%	2.333%
10	7.540	767	774	788	rBV	334276	570631	18.22%	2.167%
11	8.257	886	896	914	rBV	258042	547689	17.49%	2.080%
12	8.681	961	968	990	rBV	356169	702770	22.44%	2.669%
13	9.069	1029	1034	1040	rVB3	6794	10856	0.35%	0.041%
14	9.393	1081	1089	1107	rBV	308750	621174	19.83%	2.359%
15	9.940	1175	1182	1209	rBV	351738	894933	28.57%	3.399%
16	10.281	1233	1240	1251	rBV	443134	752609	24.03%	2.858%
17	10.446	1260	1268	1295	rBV	244996	596537	19.05%	2.266%
18	11.528	1448	1452	1457	rBV6	2937	5042	0.16%	0.019%
19	11.916	1511	1518	1526	rBV7	3946	10557	0.34%	0.040%
20	13.557	1790	1797	1814	rBV	965153	1429347	45.64%	5.428%
21	13.839	1837	1845	1862	rBV	1043527	1480580	47.27%	5.623%
22	14.157	1890	1899	1908	rBV	706201	1068603	34.12%	4.058%
23	14.492	1950	1956	1968	rBV5	28709	105998	3.38%	0.403%
24	15.169	2063	2071	2082	rBV	1370265	1964155	62.71%	7.459%
25	15.322	2091	2097	2111	rBV	308998	626879	20.02%	2.381%
26	15.422	2112	2114	2118	rVB3	10308	11440	0.37%	0.043%
27	15.545	2130	2135	2147	rBV	72915	109536	3.50%	0.416%
28	16.128	2228	2234	2246	rVB3	18062	34203	1.09%	0.130%
29	16.957	2369	2375	2386	rVV	952193	1370242	43.75%	5.204%
30	17.063	2386	2393	2413	rVB	1386709	2235243	71.37%	8.489%
31	17.898	2530	2535	2541	rBV	41457	58268	1.86%	0.221%
32	18.186	2581	2584	2594	rVB	5822	13381	0.43%	0.051%
33	18.345	2608	2611	2619	rBV9	3929	8562	0.27%	0.033%
34	18.769	2679	2683	2687	rVB4	14141	22199	0.71%	0.084%
35	18.992	2717	2721	2727	rVB3	20070	32896	1.05%	0.125%
36	19.086	2732	2737	2744	rBV	18400	34209	1.09%	0.130%

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37	19.239	2760	2763	2768	rBV5	16179	20010	0.64%	0.076%
38	19.451	2792	2799	2814	rBV	1779198	2771025	88.48%	10.524%
39	21.386	3123	3128	3145	rBV2	1034021	1695026	54.12%	6.437%
40	24.351	3621	3632	3647	rVB	1061218	3131923	100.00%	11.894%
41	24.562	3658	3668	3685	rVB	557584	1937471	61.86%	7.358%

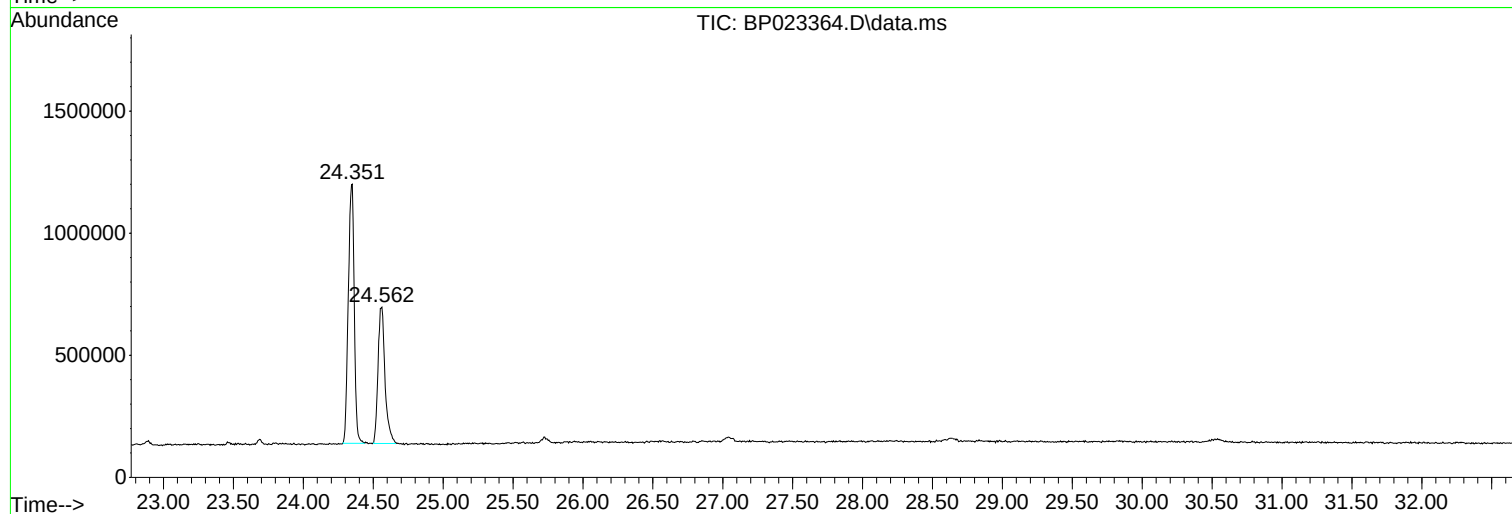
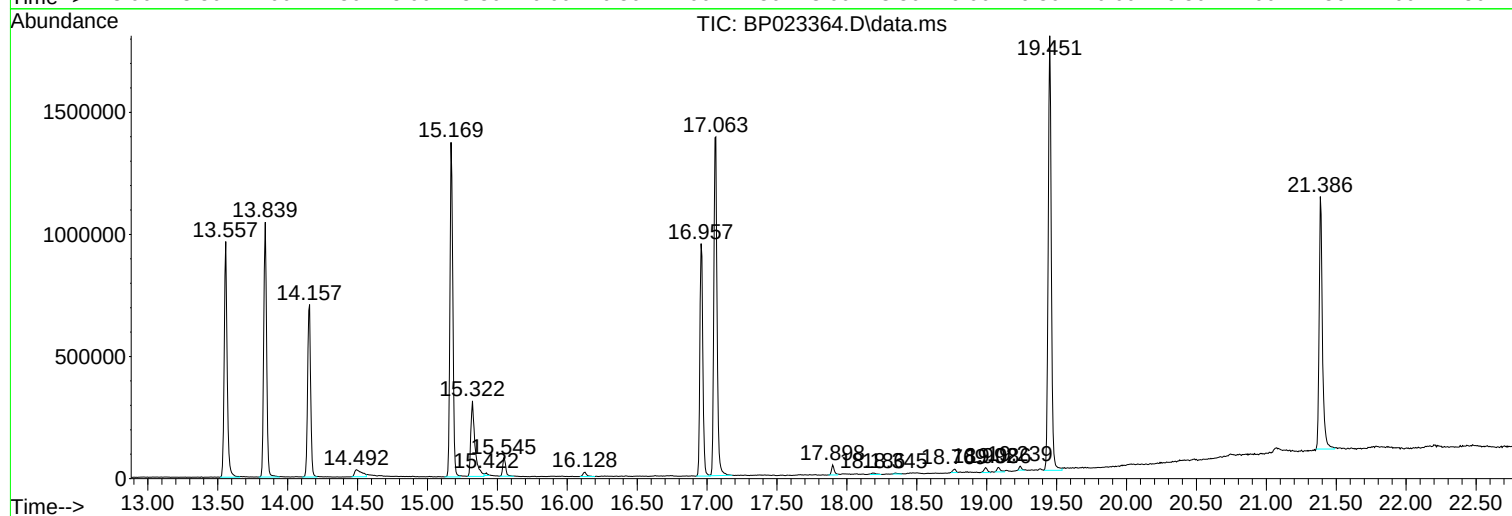
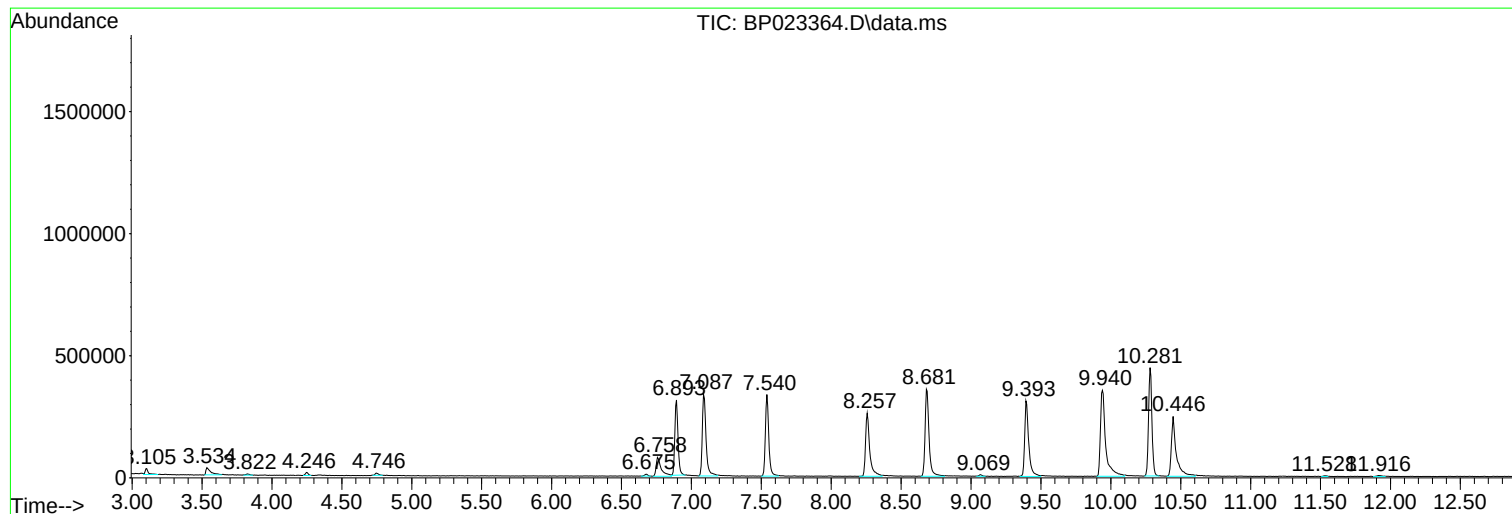
Sum of corrected areas: 26331351

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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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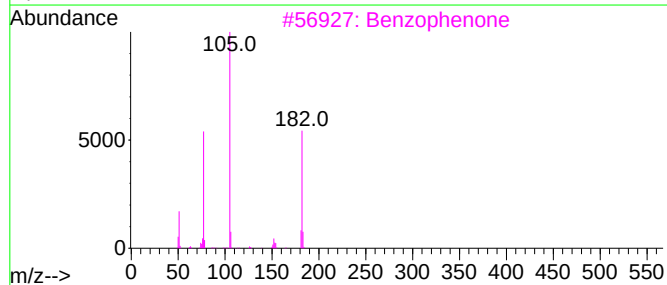
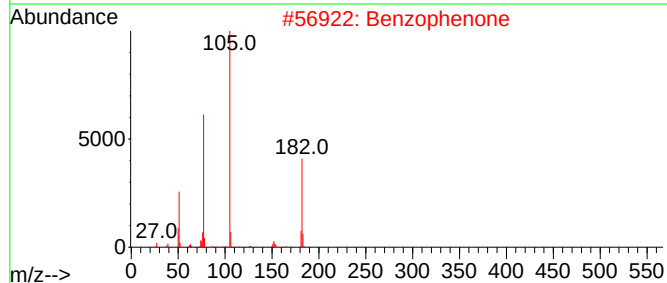
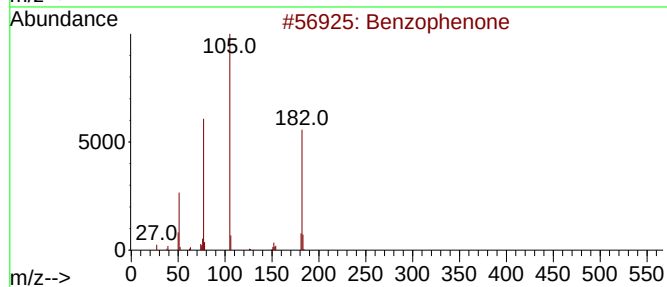
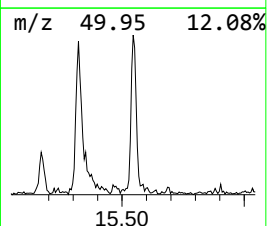
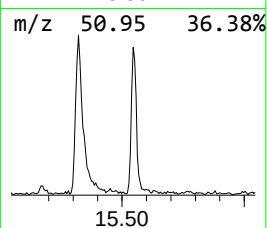
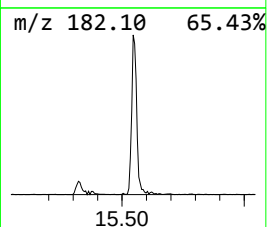
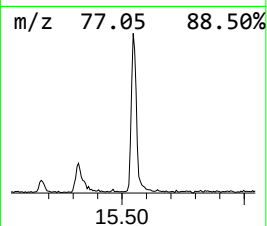
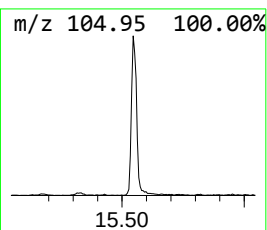
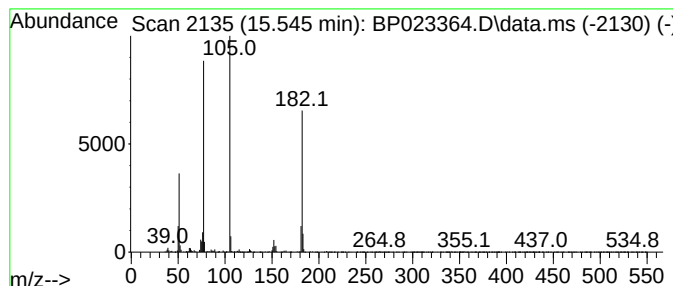
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
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TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Benzophenone Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.545	2.05 ng/ul	109536	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	95
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 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	2.0	ng/u1	109536	3	14.157	1068600	20.0