

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050823\
 Data File : BP015010.D
 Acq On : 09 May 2023 01:10
 Operator : CG/JU
 Sample : 02609-08
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 JREH3

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-BP050323.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP015010.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.458	8	12	18	rVB	75639	104082	0.85%	0.108%
2	3.899	84	87	104	rBV	279759	460969	3.76%	0.478%
3	4.140	124	128	132	rBV4	12288	18694	0.15%	0.019%
4	4.240	142	145	152	rVB	25602	37397	0.30%	0.039%
5	7.305	655	666	677	rBV	513579	829367	6.76%	0.861%
6	7.475	688	695	708	rBV	1249240	1951911	15.91%	2.026%
7	7.681	723	730	741	rBV	1181024	1914783	15.61%	1.987%
8	8.157	804	811	820	rBV	1024869	1644575	13.41%	1.707%
9	8.869	924	932	941	rBV	1196832	1942293	15.83%	2.016%
10	9.340	1004	1012	1025	rBV	1457358	2480475	20.22%	2.574%
11	10.063	1127	1135	1148	rBV	1021122	1739240	14.18%	1.805%
12	10.610	1219	1228	1243	rBV	1692268	2964241	24.16%	3.077%
13	10.993	1284	1293	1304	rBV	1568046	2694685	21.97%	2.797%
14	11.128	1309	1316	1334	rBV	990292	1677529	13.68%	1.741%
15	12.234	1497	1504	1511	rVB	70742	106393	0.87%	0.110%
16	12.569	1554	1561	1569	rBV2	27211	52262	0.43%	0.054%
17	13.687	1746	1751	1759	rVV2	33788	54605	0.45%	0.057%
18	13.775	1761	1766	1773	rVB5	7530	15441	0.13%	0.016%
19	14.210	1832	1840	1852	rBV	3681257	5379756	43.86%	5.584%
20	14.322	1855	1859	1865	rVB8	7977	14007	0.11%	0.015%
21	14.498	1882	1889	1904	rBV	5070458	7725936	62.98%	8.019%
22	14.804	1934	1941	1953	rBV	3123608	4627224	37.72%	4.803%
23	14.986	1966	1972	1992	rBV	369922	682393	5.56%	0.708%
24	15.210	2006	2010	2018	rVB10	8944	18436	0.15%	0.019%
25	15.798	2103	2110	2123	rVB	6477085	9353444	76.25%	9.708%
26	15.910	2123	2129	2140	rVB	352836	511918	4.17%	0.531%
27	16.033	2146	2150	2158	rVB2	34473	56710	0.46%	0.059%
28	16.157	2165	2171	2180	rBV	367872	514079	4.19%	0.534%
29	16.498	2225	2229	2235	rBV9	6963	13599	0.11%	0.014%
30	17.233	2350	2354	2360	rVB7	8063	13144	0.11%	0.014%
31	17.345	2370	2373	2380	rVB4	10502	17351	0.14%	0.018%
32	17.563	2403	2410	2420	rBV	3784313	5473111	44.62%	5.680%
33	17.663	2420	2427	2438	rBV	7653708	11012837	89.78%	11.430%
34	19.457	2727	2732	2737	rBV2	106374	141148	1.15%	0.146%
35	19.522	2739	2743	2749	rVV	85036	124373	1.01%	0.129%
36	19.780	2784	2787	2791	rBV2	29566	33019	0.27%	0.034%

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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-BP050323.MA.M
Title : SVOA CALIBRATION

37	19.880	2798	2804	2819	rBV	9156436	12266831	100.00%	12.732%
38	21.310	3045	3047	3052	rBV2	85788	99916	0.81%	0.104%
39	21.645	3098	3104	3113	rBV	3523151	4843640	39.49%	5.027%
40	24.074	3509	3517	3530	rBV2	4009951	8852804	72.17%	9.188%
41	24.245	3539	3546	3555	rVB2	1809391	3884640	31.67%	4.032%

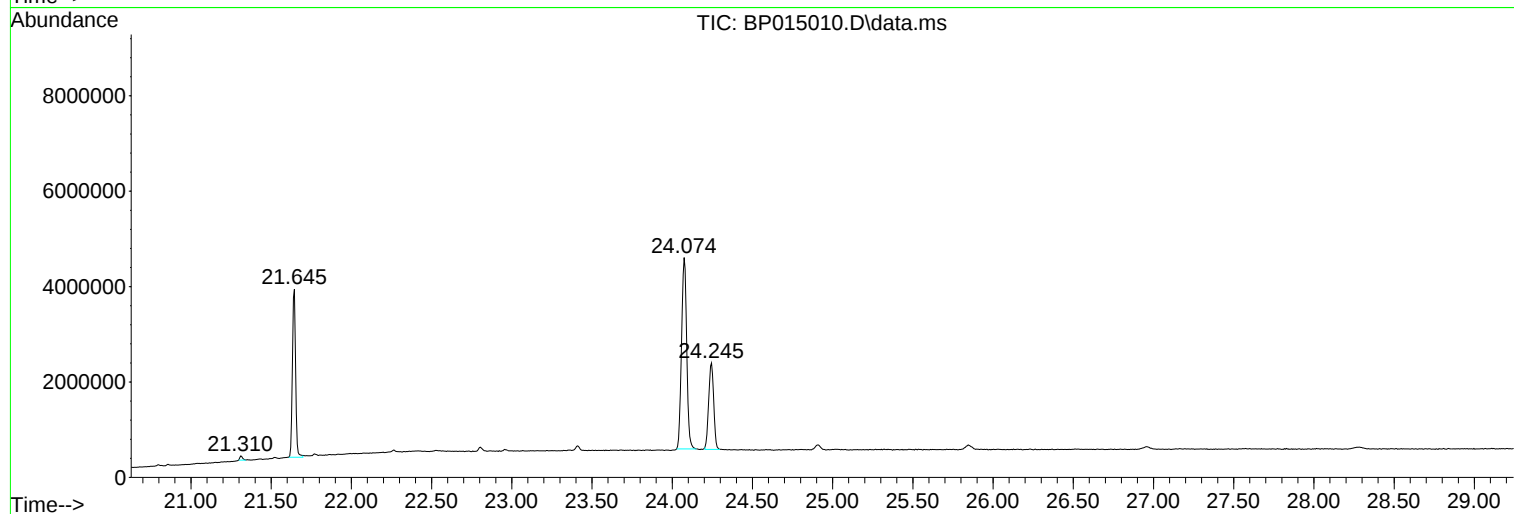
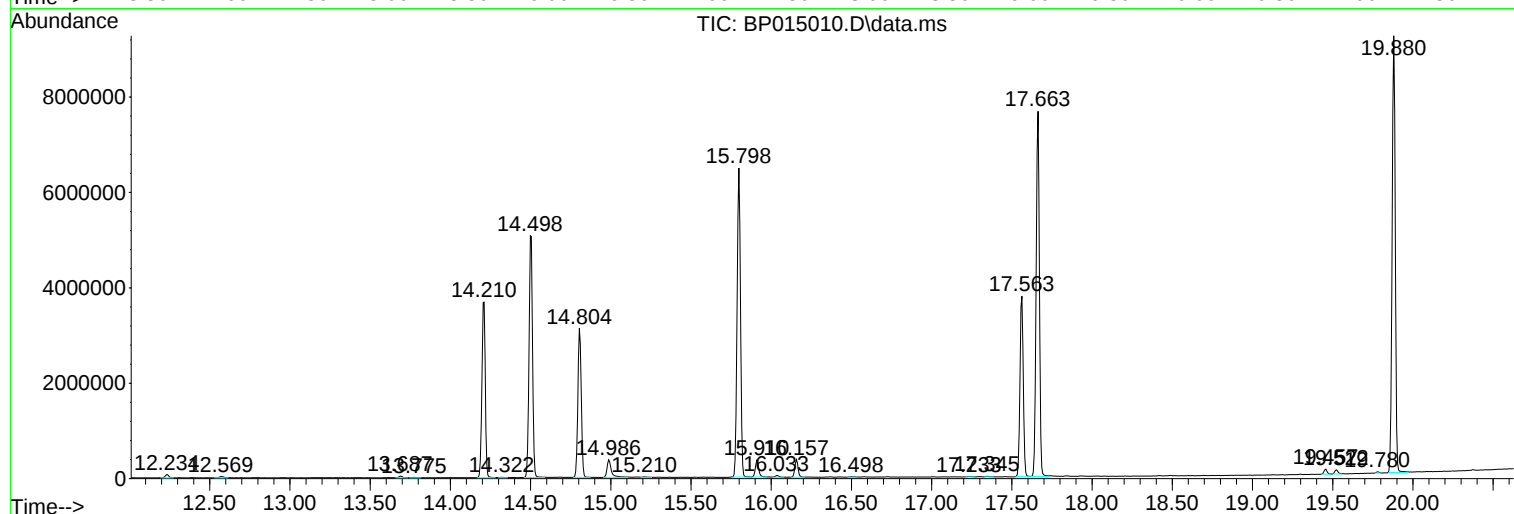
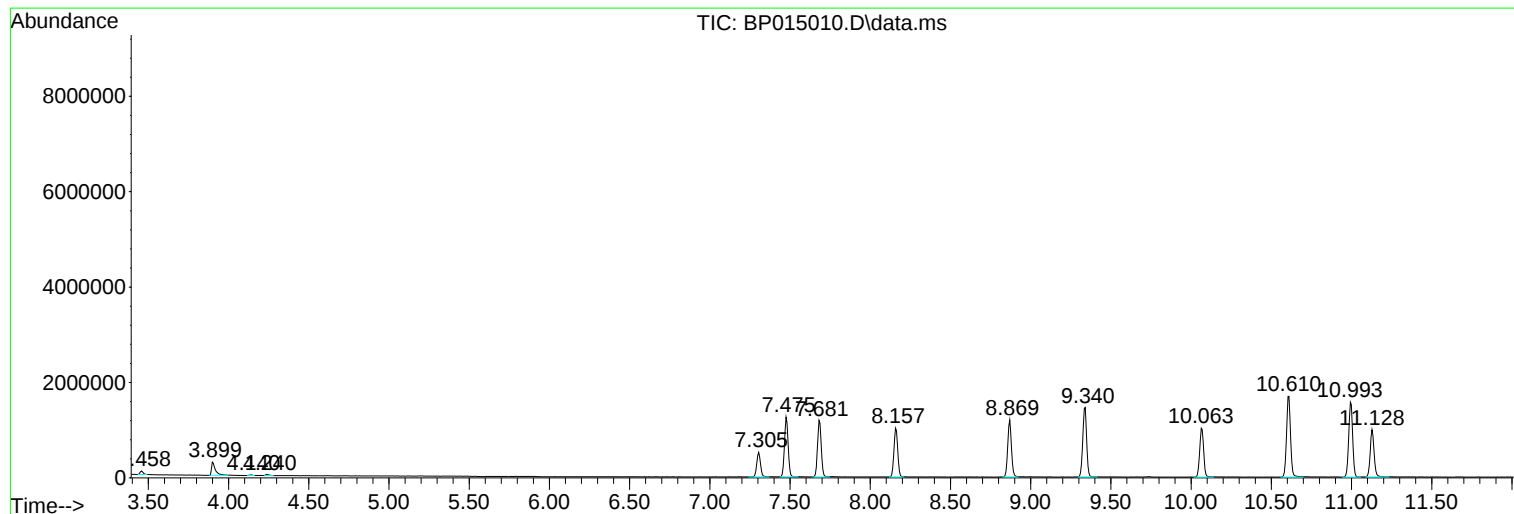
Sum of corrected areas: 96349258

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050323.MA.M
Quant Title : SVOA CALIBRATION

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



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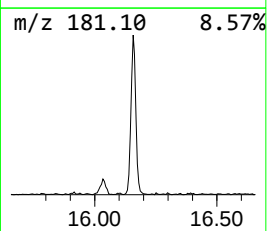
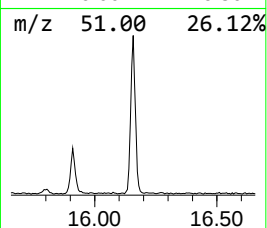
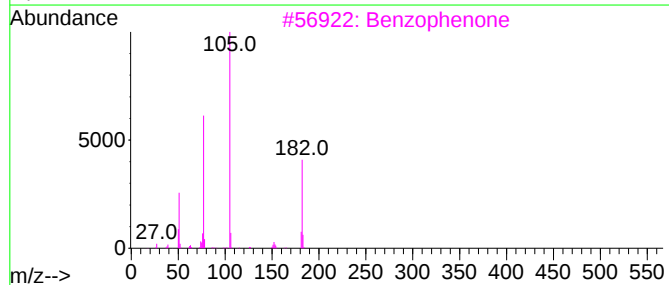
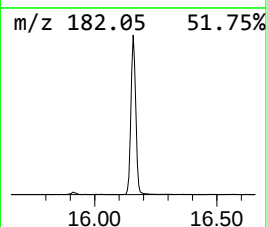
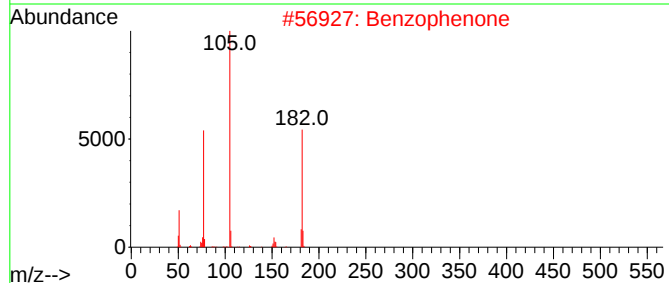
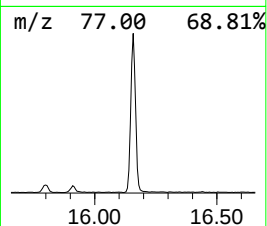
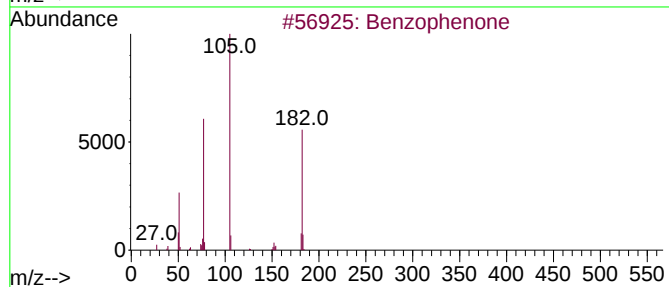
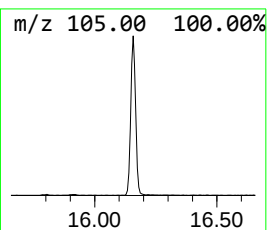
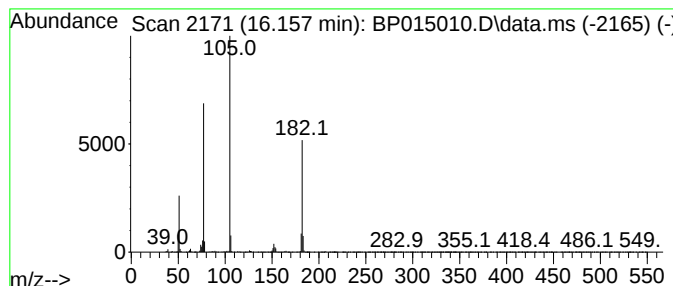
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050323.MA.M
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TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Benzophenone Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.157	2.22 ng/ul	514079	Acenaphthene-d10	14.804

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	94
4			Benzophenone	182	C13H10O	000119-61-9	93
5			Methanone, phenyl-2-pyridinyl-	183	C12H9NO	000091-02-1	72



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Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzophenone	16.157	2.2	ng/u1	514079	3	14.804	4627220	20.0