

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061423\
 Data File : BP015532.D
 Acq On : 14 Jun 2023 23:22
 Operator : MA/JU
 Sample : 03104-04
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 CONM0

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-BP060723.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP015532.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.369	26	31	36	rBV	218870	342034	5.49%	0.679%
2	3.857	112	114	129	rBV	138358	271893	4.37%	0.540%
3	4.081	147	152	156	rBV2	12143	18991	0.30%	0.038%
4	4.181	167	169	174	rVB4	5670	7260	0.12%	0.014%
5	5.152	331	334	341	rVB2	6690	10168	0.16%	0.020%
6	7.246	683	690	702	rBV	158533	289048	4.64%	0.574%
7	7.410	711	718	730	rBV	584757	941247	15.11%	1.870%
8	7.610	743	752	764	rBV	575541	1008028	16.18%	2.003%
9	8.087	827	833	845	rBV	417123	692925	11.12%	1.377%
10	8.798	944	954	968	rBV	505694	904249	14.52%	1.796%
11	9.263	1025	1033	1050	rBV	803089	1375763	22.09%	2.733%
12	9.645	1095	1098	1104	rVB7	4109	7021	0.11%	0.014%
13	9.992	1148	1157	1174	rBV	701593	1201463	19.29%	2.387%
14	10.534	1242	1249	1272	rBV	846400	1628649	26.15%	3.235%
15	10.916	1305	1314	1326	rVB	665817	1130810	18.15%	2.246%
16	11.057	1329	1338	1357	rBV	627489	1320100	21.19%	2.622%
17	12.510	1580	1585	1595	rBV3	13765	30154	0.48%	0.060%
18	13.610	1767	1772	1778	rBV3	21982	35173	0.56%	0.070%
19	14.139	1853	1862	1876	rBV	2109573	3018669	48.46%	5.997%
20	14.239	1878	1879	1887	rVB8	4758	8099	0.13%	0.016%
21	14.427	1904	1911	1929	rBV2	2200024	3324727	53.38%	6.605%
22	14.733	1956	1963	1980	rVB	1115856	1678076	26.94%	3.334%
23	14.969	1997	2003	2023	rBV5	44815	196458	3.15%	0.390%
24	15.727	2125	2132	2146	rBV	2859826	4359406	69.99%	8.660%
25	15.851	2147	2153	2167	rVV	1075014	1621715	26.04%	3.222%
26	15.969	2169	2173	2179	rVB3	19285	33751	0.54%	0.067%
27	16.086	2187	2193	2200	rBV	129052	193677	3.11%	0.385%
28	16.657	2286	2290	2296	rVB4	10095	15548	0.25%	0.031%
29	17.163	2371	2376	2380	rBV8	5722	8781	0.14%	0.017%
30	17.486	2424	2431	2442	rBV	1487733	2139671	34.35%	4.251%
31	17.592	2442	2449	2465	rVV	3393361	5157821	82.81%	10.246%
32	18.345	2573	2577	2586	rBV2	59195	96482	1.55%	0.192%
33	19.392	2750	2755	2760	rBV	48386	70200	1.13%	0.139%
34	19.463	2762	2767	2772	rVV	47817	72958	1.17%	0.145%
35	19.515	2773	2776	2780	rVV4	14945	18505	0.30%	0.037%
36	19.568	2782	2785	2792	rVV3	33865	50434	0.81%	0.100%

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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

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37	19.810	2820	2826	2839	rBV	4614623	6228844	100.00%	12.374%
38	21.568	3120	3125	3136	rBV	1700891	2434854	39.09%	4.837%
39	23.951	3522	3530	3545	rBV2	2902261	6117691	98.22%	12.153%
40	24.115	3551	3558	3570	rVB	1034629	2276515	36.55%	4.522%

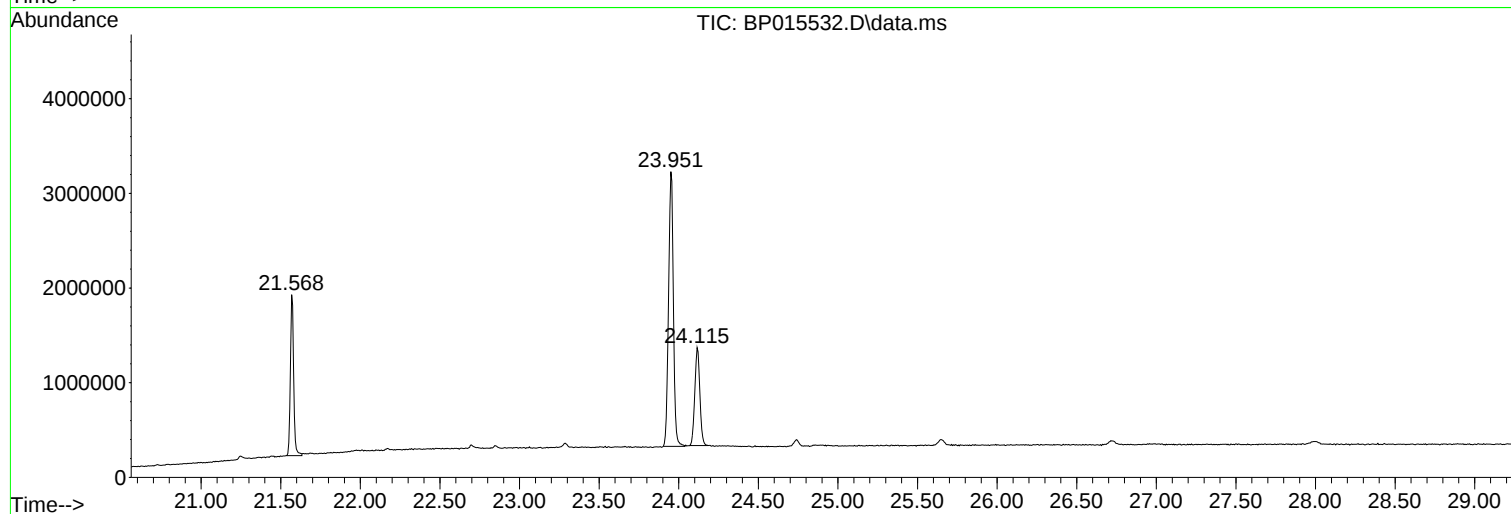
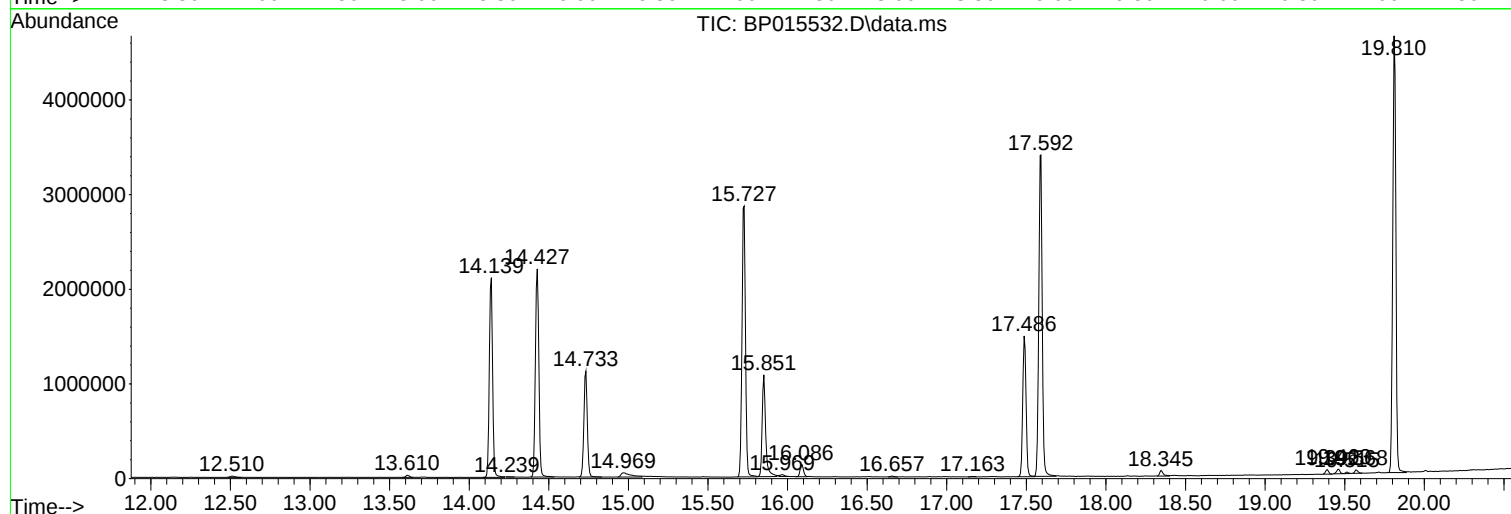
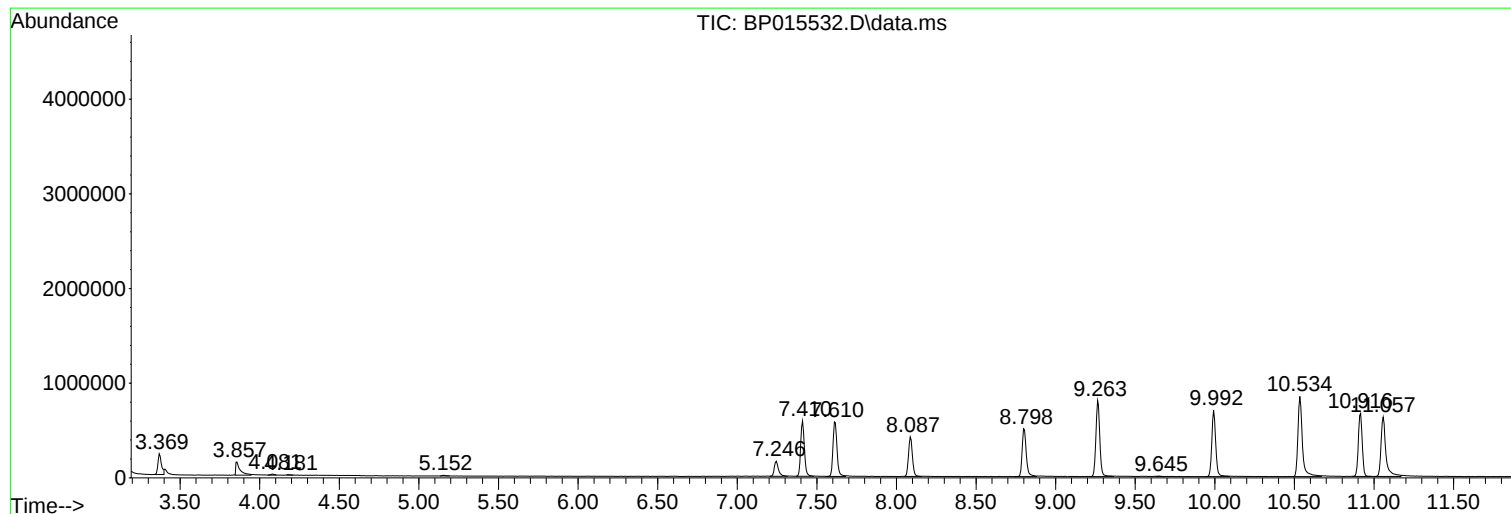
Sum of corrected areas: 50337858

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

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



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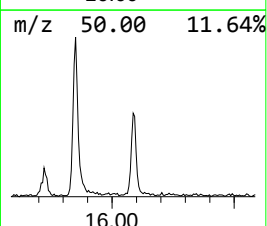
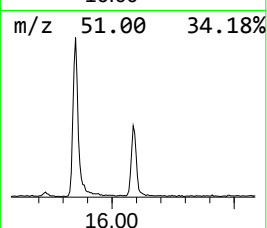
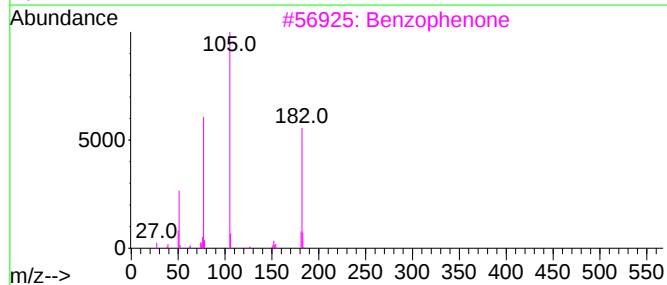
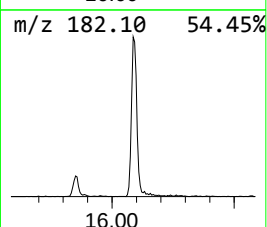
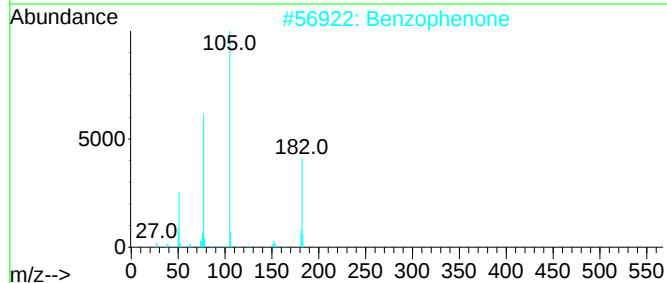
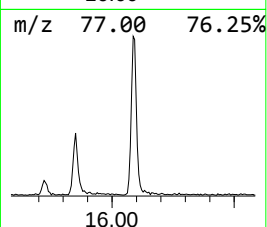
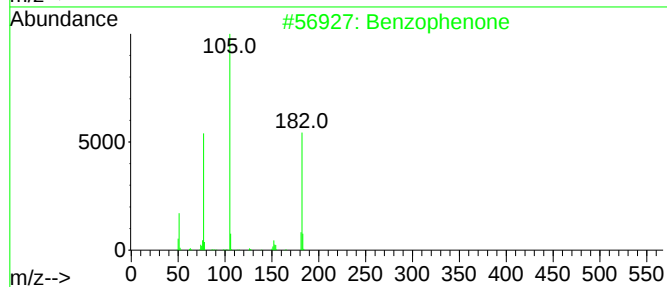
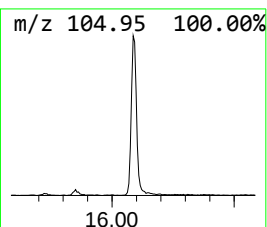
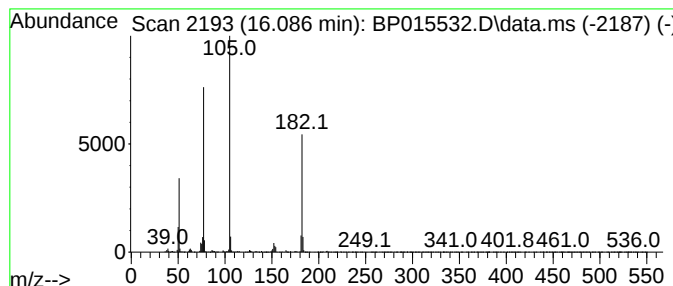
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
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TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Benzophenone Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.086	2.31 ng/ul	193677	Acenaphthene-d10	14.733

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	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\NIST20.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzophenone	16.086	2.3	ng/u1	193677	3	14.733	1678080	20.0