

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070623\
 Data File : BP016041.D
 Acq On : 06 Jul 2023 20:20
 Operator : MA/JU
 Sample : 03257-08
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EXYY0

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-BP062623.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP016041.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.352	25	28	38	rVB	51433	78712	0.98%	0.125%
2	3.817	104	107	123	rBV	182677	406982	5.04%	0.644%
3	4.017	137	141	145	rVB3	12197	19089	0.24%	0.030%
4	4.117	155	158	166	rVB4	13285	25024	0.31%	0.040%
5	7.258	684	692	704	rBV2	119402	380319	4.71%	0.602%
6	7.363	704	710	736	rVV	482560	1326361	16.44%	2.099%
7	7.581	740	747	769	rVV	435875	1214115	15.05%	1.922%
8	8.034	817	824	844	rBV	396998	1129197	14.00%	1.787%
9	8.793	944	953	980	rBV	308294	1078686	13.37%	1.707%
10	9.234	1021	1028	1054	rBV	667252	1670863	20.71%	2.645%
11	9.546	1080	1081	1090	rVB9	8871	13382	0.17%	0.021%
12	9.963	1145	1152	1175	rBV	397506	1358321	16.84%	2.150%
13	10.546	1243	1251	1273	rBV2	353945	1597217	19.80%	2.528%
14	10.863	1297	1305	1318	rBV	713988	1701085	21.08%	2.692%
15	11.051	1329	1337	1371	rBV2	422898	1811664	22.45%	2.867%
16	12.510	1581	1585	1592	rBV5	11079	24085	0.30%	0.038%
17	13.563	1759	1764	1769	rBV4	7135	11964	0.15%	0.019%
18	14.098	1847	1855	1881	rBV	1753571	3750436	46.48%	5.936%
19	14.381	1895	1903	1923	rBV2	2298427	4121188	51.08%	6.523%
20	14.681	1948	1954	1976	rBV	1392361	2411610	29.89%	3.817%
21	15.087	2013	2023	2024	rBV5	36971	80851	1.00%	0.128%
22	15.451	2081	2085	2091	rVB3	10628	17677	0.22%	0.028%
23	15.687	2118	2125	2143	rBV	2536098	5391866	66.83%	8.534%
24	15.869	2150	2156	2183	rBV	479693	1646836	20.41%	2.607%
25	16.063	2184	2189	2207	rVB	193136	376059	4.66%	0.595%
26	16.616	2278	2283	2292	rBV2	27381	50390	0.62%	0.080%
27	16.839	2316	2321	2324	rBV7	5732	8813	0.11%	0.014%
28	17.445	2418	2424	2435	rBV	1924605	2964439	36.74%	4.692%
29	17.551	2435	2442	2468	rVB2	2760022	6341744	78.60%	10.038%
30	18.316	2567	2572	2580	rBV4	54278	146171	1.81%	0.231%
31	19.357	2745	2749	2755	rBV2	60047	82896	1.03%	0.131%
32	19.428	2758	2761	2769	rVV	42338	77834	0.96%	0.123%
33	19.539	2777	2780	2783	rBV4	32333	49554	0.61%	0.078%
34	19.780	2814	2821	2847	rBV	3866454	8068323	100.00%	12.770%
35	21.551	3117	3122	3138	rBV	1629125	3509034	43.49%	5.554%
36	23.910	3517	3523	3544	rVB2	2768830	7151753	88.64%	11.320%

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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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Title : SVOA CALIBRATION

37	24.086	3547	3553	3571	rVB	999073	3085016	38.24%	4.883%
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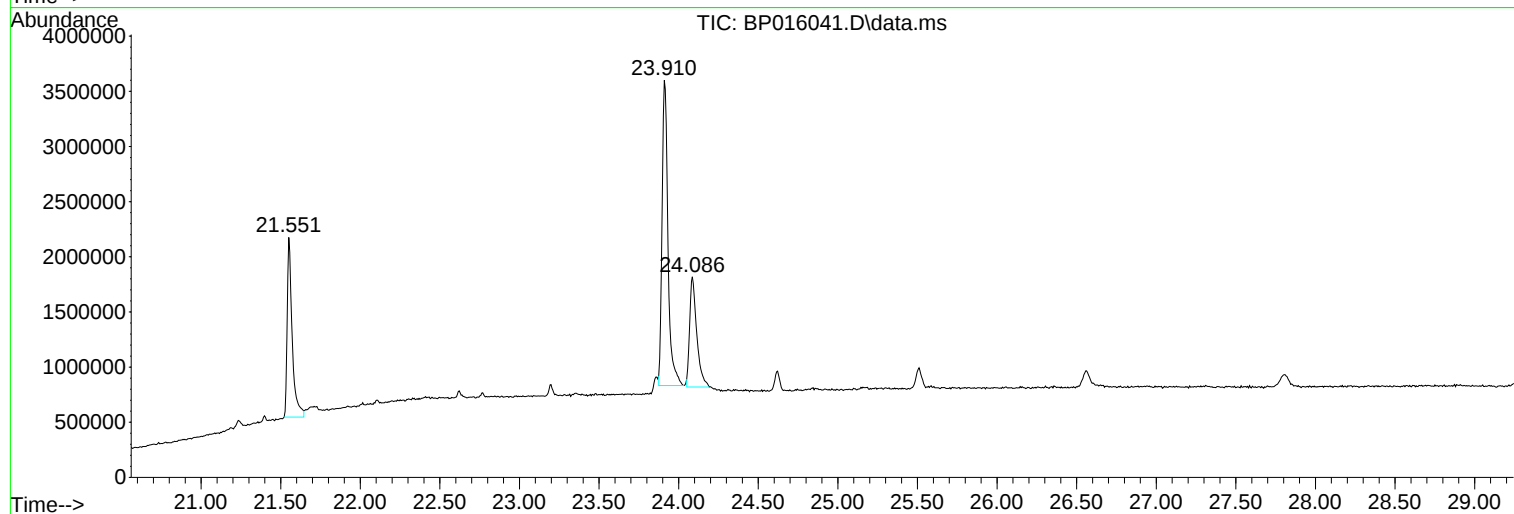
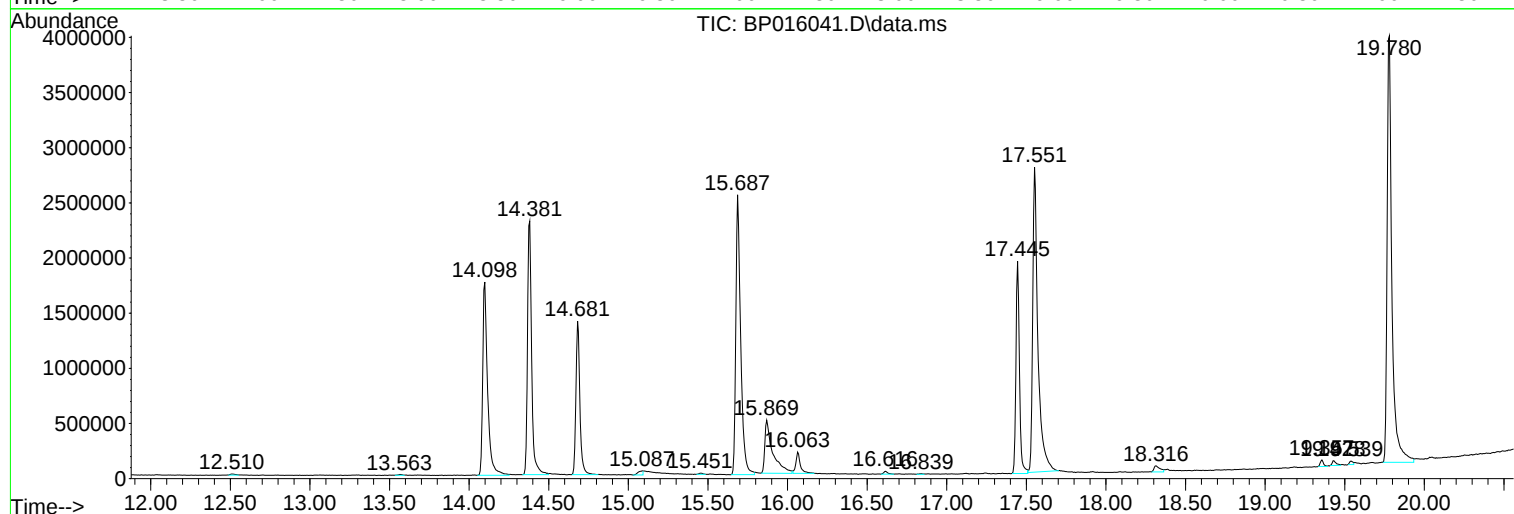
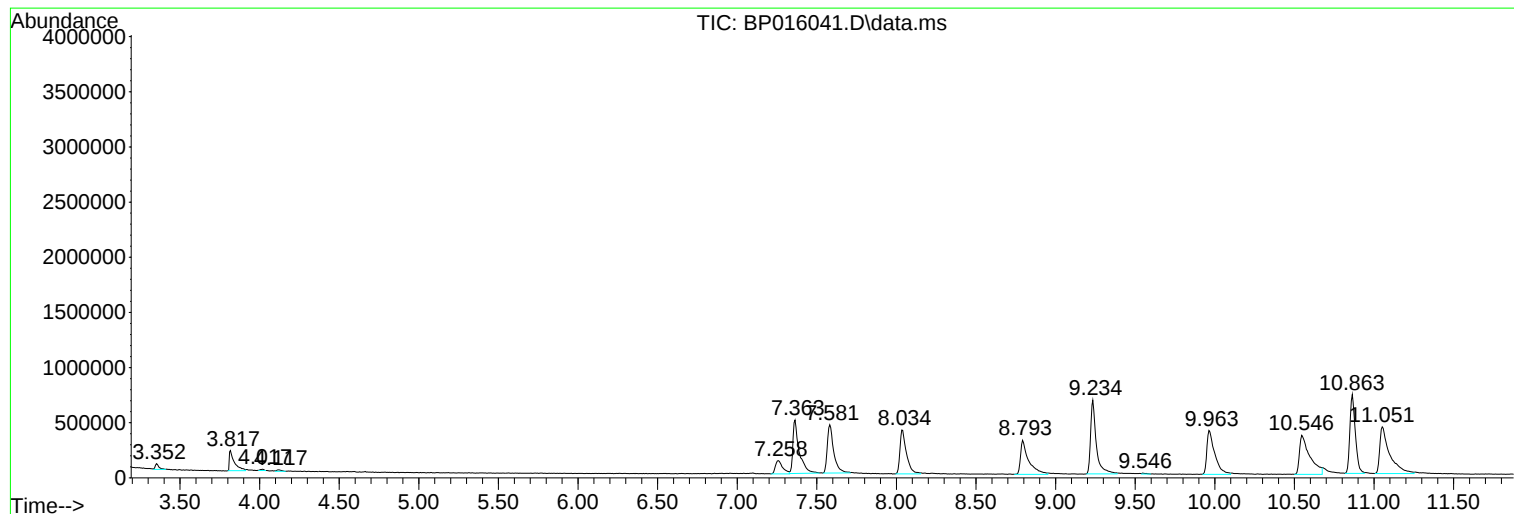
Sum of corrected areas: 63179556

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

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



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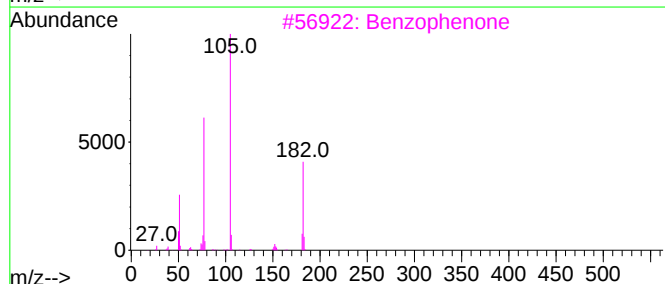
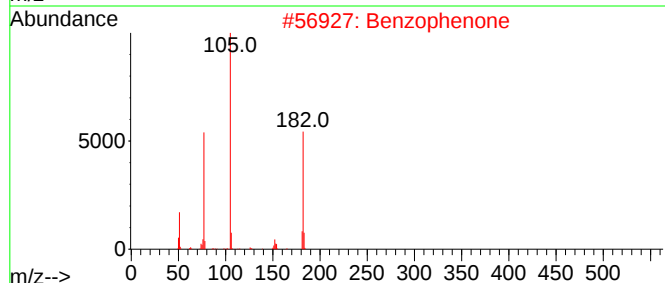
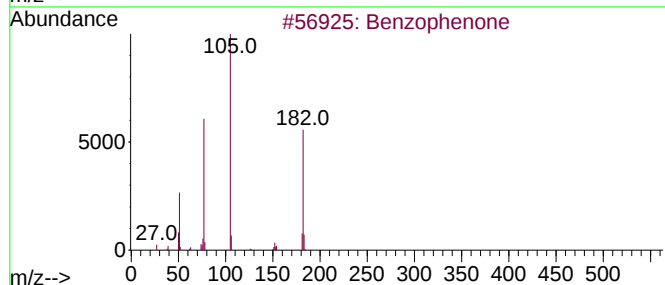
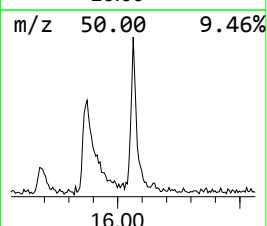
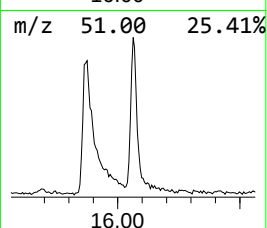
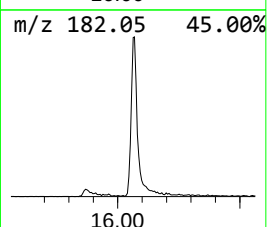
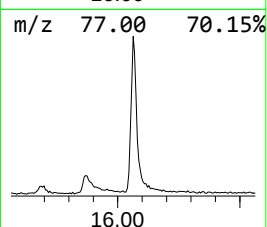
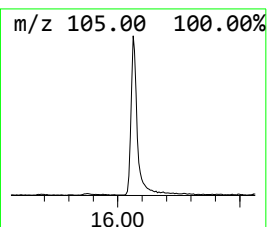
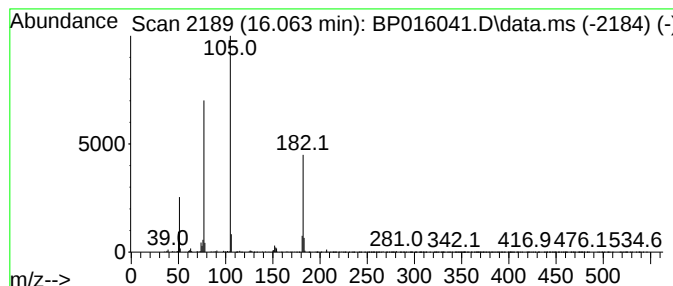
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TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Benzophenone Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.063	2.54 ng/ul	376059	Phenanthrene-d10	17.445

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	91
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	16.063	2.5	ng/u1	376059	4	17.445	2964440	20.0