

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042823\
 Data File : BP014892.D
 Acq On : 29 Apr 2023 01:30
 Operator : CG/JU
 Sample : 02528-02
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 HOA34

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-BP042823.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP014892.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.434	3	8	16	rBV	1719587	2552762	8.87%	3.107%
2	3.505	16	20	33	rVB	1063112	1420668	4.94%	1.729%
3	3.911	86	89	108	rBV	203699	356214	1.24%	0.434%
4	4.152	126	130	137	rVB	22843	33219	0.12%	0.040%
5	4.811	235	242	266	rVB	14959352	28777439	100.00%	35.031%
6	5.828	406	415	422	rBV	17238	31729	0.11%	0.039%
7	7.311	660	667	682	rBV	146612	267193	0.93%	0.325%
8	7.487	690	697	715	rBV	754841	1236568	4.30%	1.505%
9	7.693	724	732	744	rBV	637457	1096010	3.81%	1.334%
10	8.175	806	814	827	rBV	578391	969627	3.37%	1.180%
11	8.875	926	933	951	rBV	440508	816040	2.84%	0.993%
12	9.352	1005	1014	1027	rBV	756784	1363489	4.74%	1.660%
13	10.081	1130	1138	1154	rBV	558344	1002942	3.49%	1.221%
14	10.616	1222	1229	1246	rBV	846238	1554978	5.40%	1.893%
15	11.010	1288	1296	1310	rBV	778316	1358567	4.72%	1.654%
16	11.140	1311	1318	1343	rBV	695346	1455295	5.06%	1.772%
17	13.704	1748	1754	1762	rBV2	21541	34166	0.12%	0.042%
18	14.216	1834	1841	1859	rBV	2019997	2886226	10.03%	3.513%
19	14.510	1884	1891	1912	rBV	2218521	3499945	12.16%	4.260%
20	14.816	1936	1943	1963	rBV	1120918	1784068	6.20%	2.172%
21	14.993	1968	1973	1981	rBV3	31984	76648	0.27%	0.093%
22	15.810	2104	2112	2125	rBV	2847565	4387788	15.25%	5.341%
23	15.916	2125	2130	2145	rVB	747969	1079835	3.75%	1.314%
24	16.169	2168	2173	2186	rBV	297666	449296	1.56%	0.547%
25	16.734	2265	2269	2275	rBV	50555	74562	0.26%	0.091%
26	17.569	2405	2411	2422	rBV	1249755	1995097	6.93%	2.429%
27	17.669	2422	2428	2457	rVB	3172826	4938549	17.16%	6.012%
28	18.428	2553	2557	2563	rBV4	31881	48034	0.17%	0.058%
29	18.516	2566	2572	2584	rBV	259894	371303	1.29%	0.452%
30	19.469	2730	2734	2739	rBV2	38611	57120	0.20%	0.070%
31	19.533	2741	2745	2750	rBV	34787	53058	0.18%	0.065%
32	19.651	2762	2765	2771	rBV4	24481	36489	0.13%	0.044%
33	19.892	2799	2806	2815	rBV	4107760	5763156	20.03%	7.015%
34	21.651	3100	3105	3120	rBV	1236437	2107883	7.32%	2.566%
35	24.086	3511	3519	3539	rBV2	2549393	5933206	20.62%	7.222%
36	24.257	3541	3548	3562	rVB2	996920	2280188	7.92%	2.776%

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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-BP042823.MA.M
Title : SVOA CALIBRATION

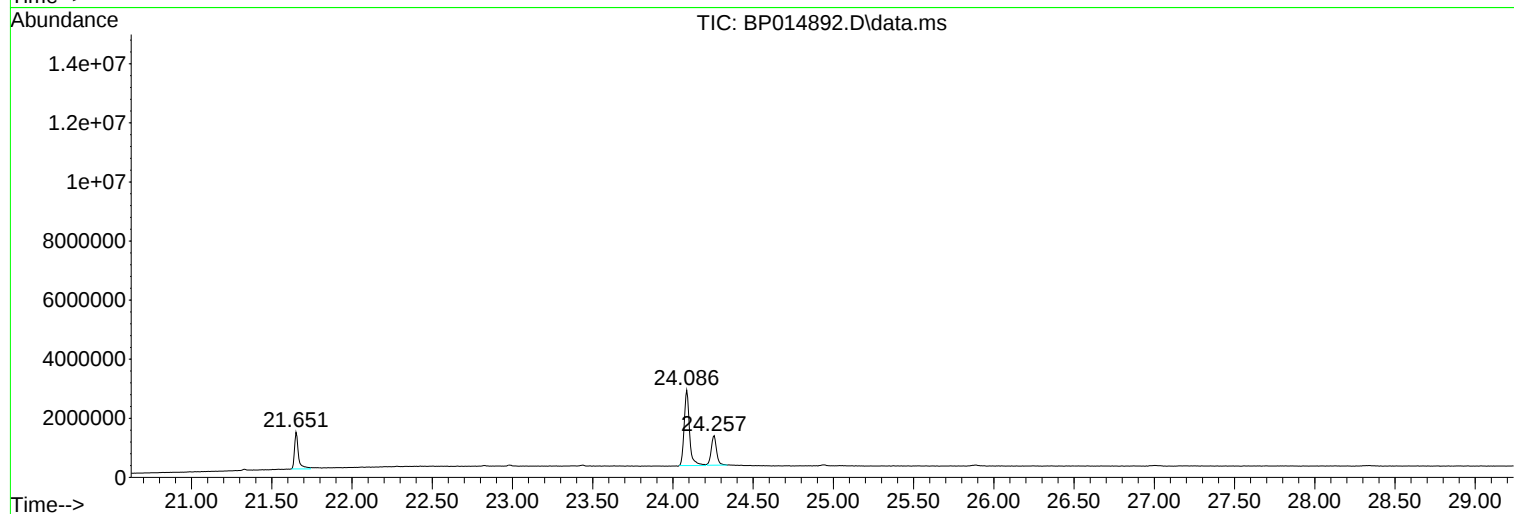
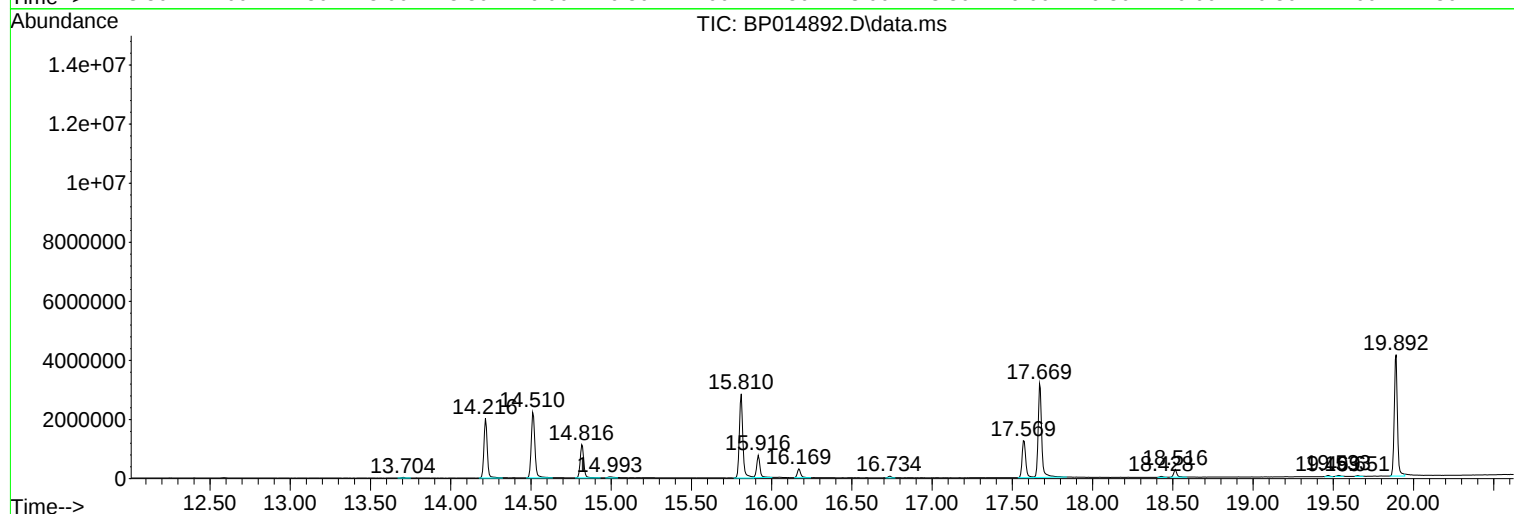
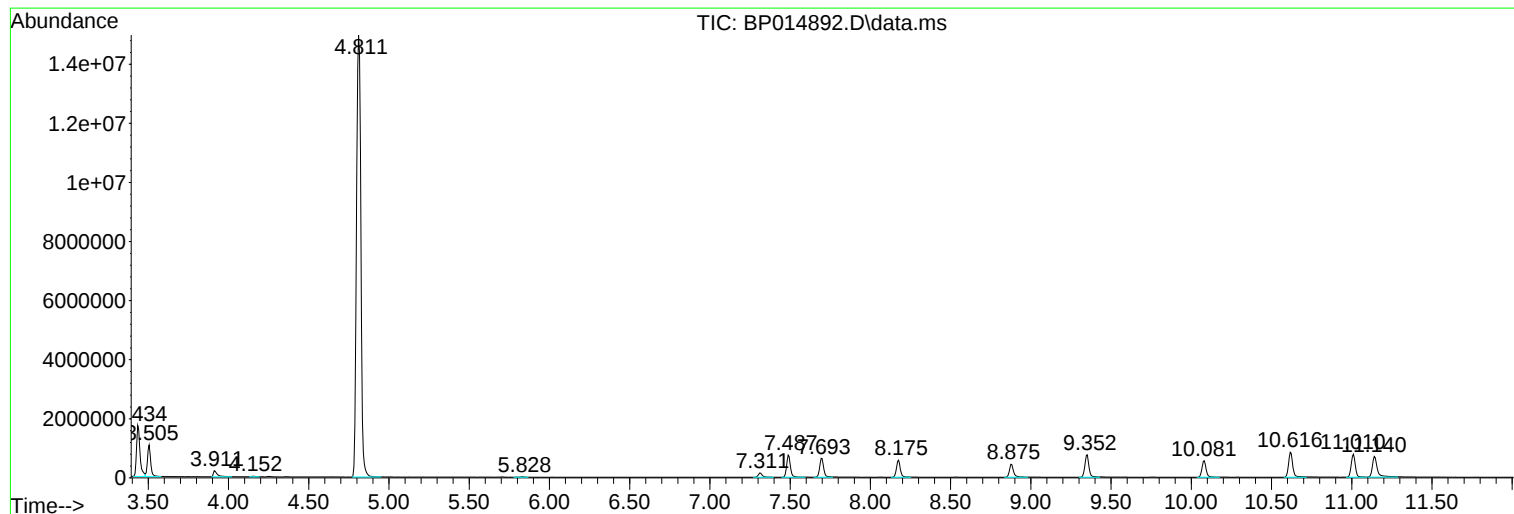
Sum of corrected areas: 82149357

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042823.MA.M
Quant Title : SVOA CALIBRATION

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



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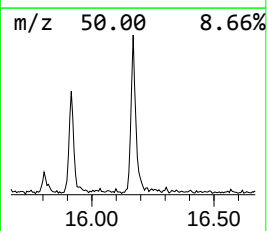
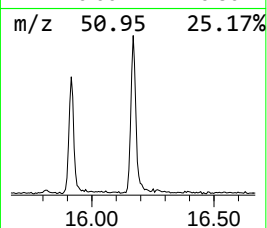
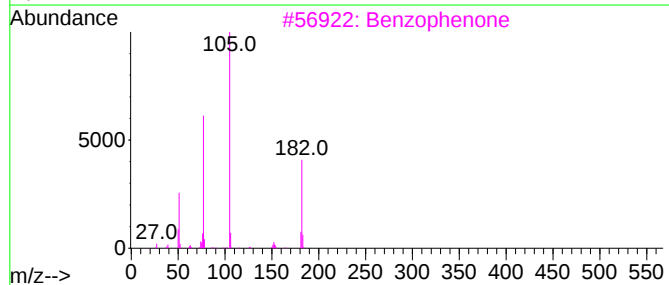
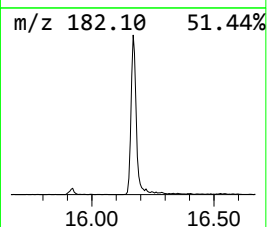
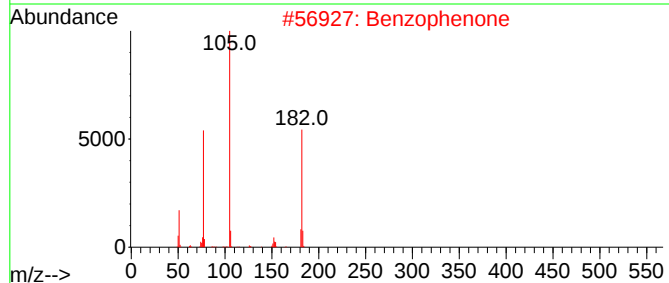
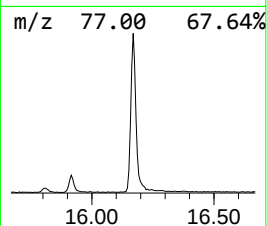
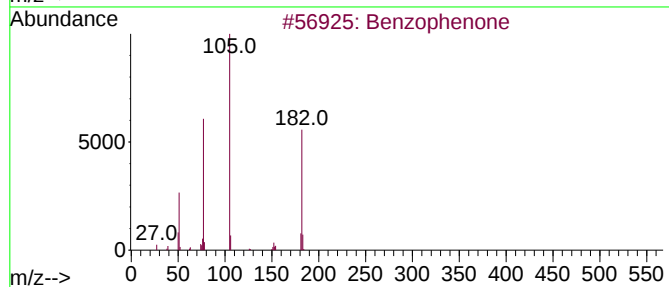
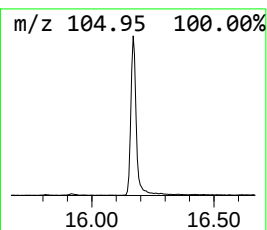
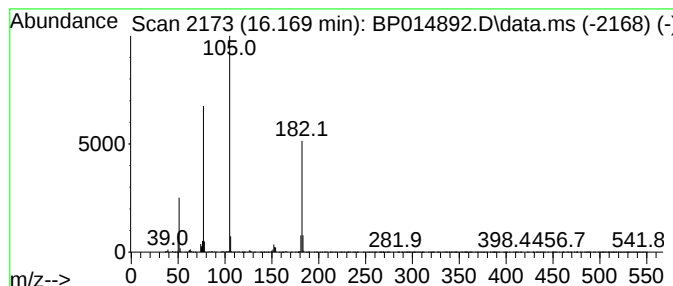
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042823.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzophenone Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.169	5.04 ng/ul	449296	Acenaphthene-d10	14.816

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			Benzophenone	182	C13H10O	000119-61-9	91



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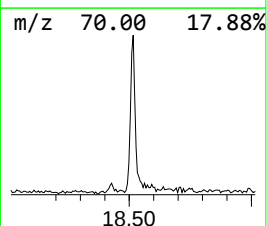
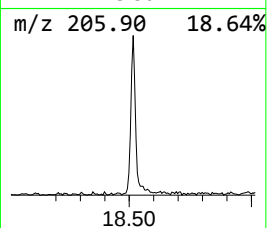
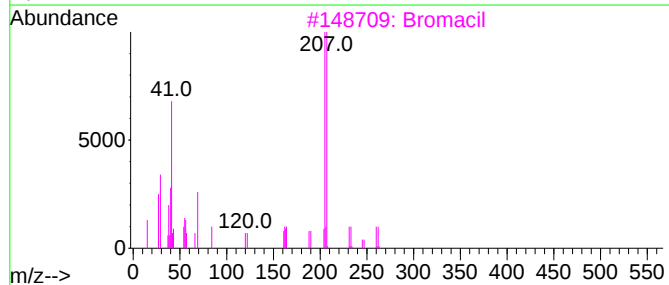
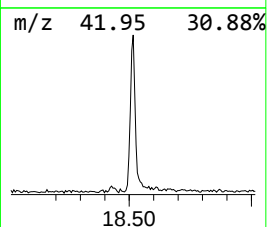
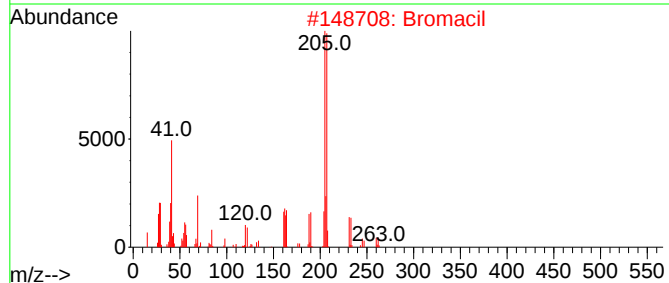
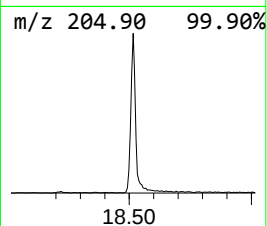
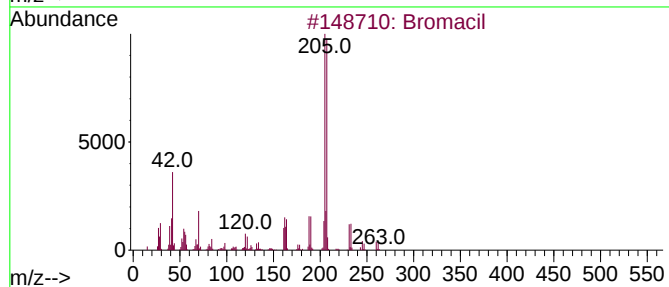
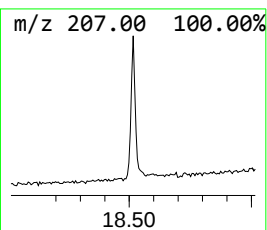
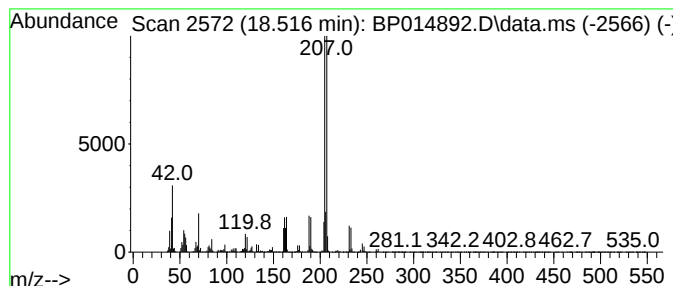
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042823.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Bromacil Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.516	3.72 ng/ul	371303	Phenanthrene-d10	17.569

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Bromacil	260	C9H13BrN2O2	000314-40-9	94
2			Bromacil	260	C9H13BrN2O2	000314-40-9	91
3			Bromacil	260	C9H13BrN2O2	000314-40-9	83
4			5-Acetamido-2,4-dichlorobenzoic ...	247	C9H7Cl2NO3	050602-49-8	64
5			Benzoic acid, 3-amino-2,5-dichloro-	205	C7H5Cl2NO2	000133-90-4	50



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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042823.MA.M
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	16.169	5.0	ng/ul	449296	3	14.816	1784070	20.0
Bromacil	18.516	3.7	ng/ul	371303	4	17.569	1995100	20.0