

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042823\
 Data File : BP014893.D
 Acq On : 29 Apr 2023 02:03
 Operator : CG/JU
 Sample : 02528-03
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 HOA35

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: BP014893.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	1988810	2853703	9.96%	3.731%
2	3.505	16	20	30	rVB	724525	954494	3.33%	1.248%
3	3.911	87	89	106	rBV	179491	301356	1.05%	0.394%
4	4.152	126	130	135	rVB	20541	30803	0.11%	0.040%
5	4.811	235	242	259	rBV	15072203	28657834	100.00%	37.471%
6	7.311	660	667	685	rBV	131324	230587	0.80%	0.301%
7	7.487	690	697	716	rBV	663051	1082660	3.78%	1.416%
8	7.693	725	732	745	rBV	567435	955507	3.33%	1.249%
9	8.175	805	814	830	rBV	689254	1193223	4.16%	1.560%
10	8.875	926	933	946	rBV	372398	688825	2.40%	0.901%
11	9.346	1005	1013	1026	rBV	646811	1159474	4.05%	1.516%
12	10.075	1130	1137	1151	rBV	468571	844512	2.95%	1.104%
13	10.616	1221	1229	1241	rBV	741826	1301147	4.54%	1.701%
14	11.010	1288	1296	1311	rBV	935242	1619827	5.65%	2.118%
15	11.140	1311	1318	1334	rBV	517107	1034627	3.61%	1.353%
16	13.704	1749	1754	1760	rVB2	21282	32568	0.11%	0.043%
17	14.216	1834	1841	1859	rBV	1740919	2457489	8.58%	3.213%
18	14.510	1885	1891	1915	rBV	1846178	2898826	10.12%	3.790%
19	14.816	1936	1943	1963	rBV	1305821	2042246	7.13%	2.670%
20	14.992	1968	1973	1985	rBV2	27224	68639	0.24%	0.090%
21	15.810	2105	2112	2124	rBV	2286814	3620147	12.63%	4.733%
22	15.916	2124	2130	2143	rVB	552688	814925	2.84%	1.066%
23	16.169	2167	2173	2189	rBV	557739	789380	2.75%	1.032%
24	16.739	2265	2270	2276	rBV	86036	127875	0.45%	0.167%
25	17.569	2402	2411	2421	rBV	1389459	2201566	7.68%	2.879%
26	17.669	2422	2428	2454	rVB	2603433	4041823	14.10%	5.285%
27	18.428	2552	2557	2566	rBV3	61324	95215	0.33%	0.124%
28	18.516	2566	2572	2582	rBV	224508	335956	1.17%	0.439%
29	19.469	2730	2734	2739	rVB2	29119	40436	0.14%	0.053%
30	19.533	2742	2745	2749	rVV	24457	31589	0.11%	0.041%
31	19.651	2761	2765	2769	rBV3	38681	51952	0.18%	0.068%
32	19.886	2799	2805	2822	rBV	3449730	4701585	16.41%	6.147%
33	21.327	3047	3050	3055	rBV3	67960	96011	0.34%	0.126%
34	21.651	3100	3105	3117	rBV	1357312	2173341	7.58%	2.842%
35	24.080	3511	3518	3534	rBV2	2017297	4570611	15.95%	5.976%
36	24.257	3541	3548	3564	rVB2	1067388	2380155	8.31%	3.112%

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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

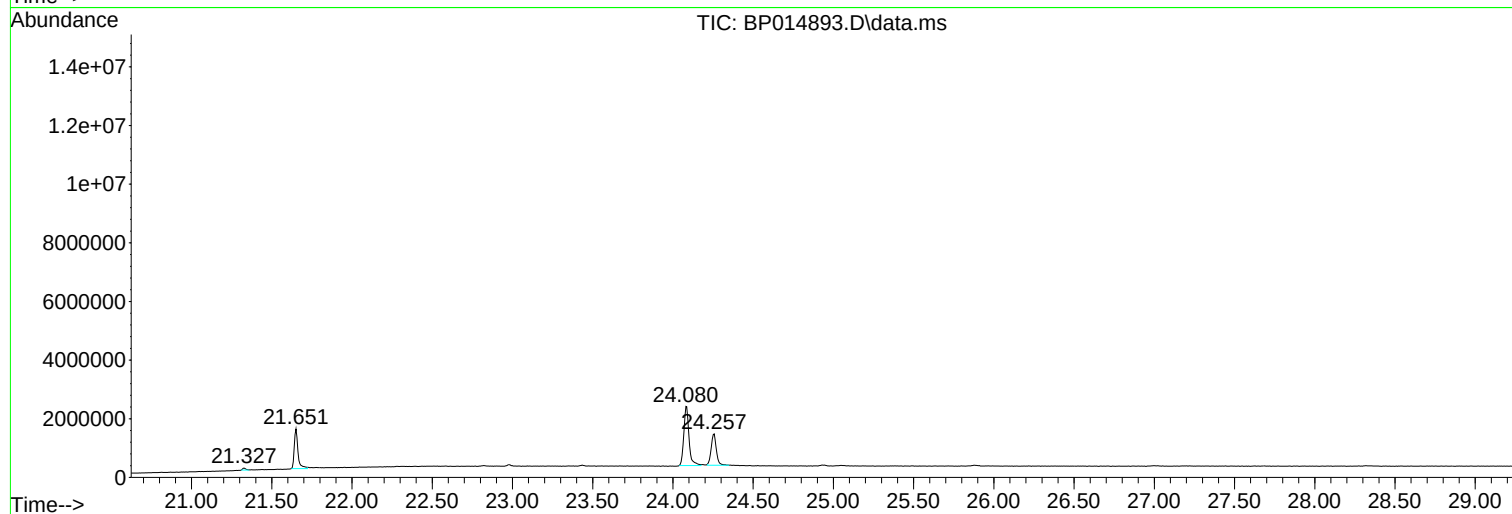
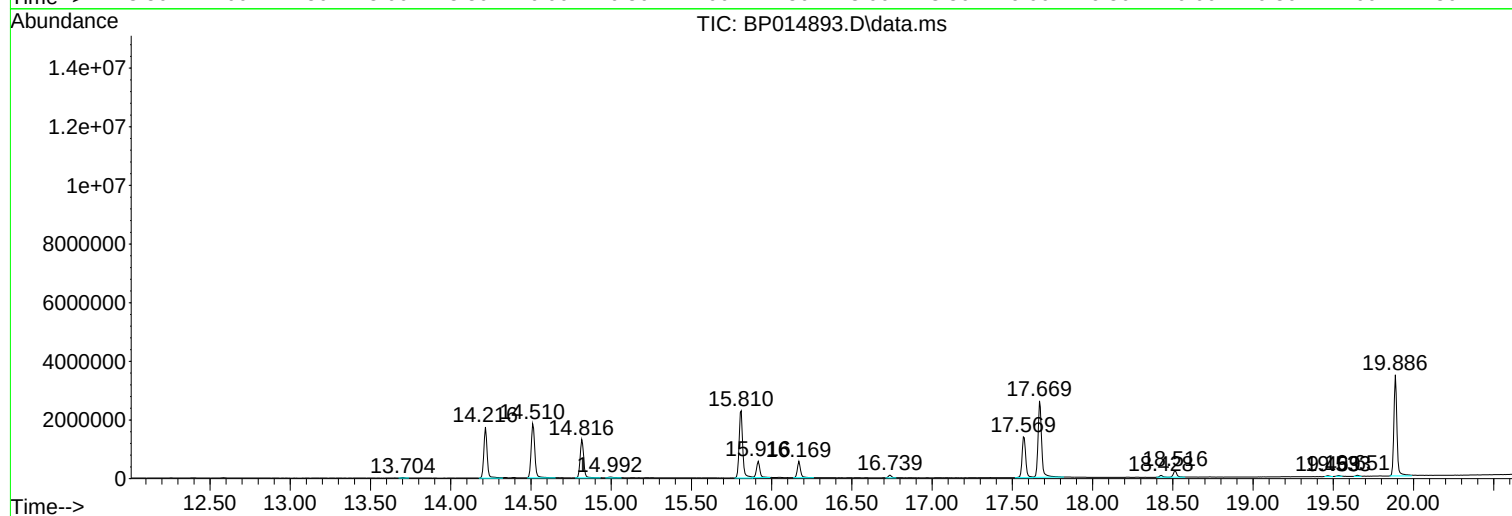
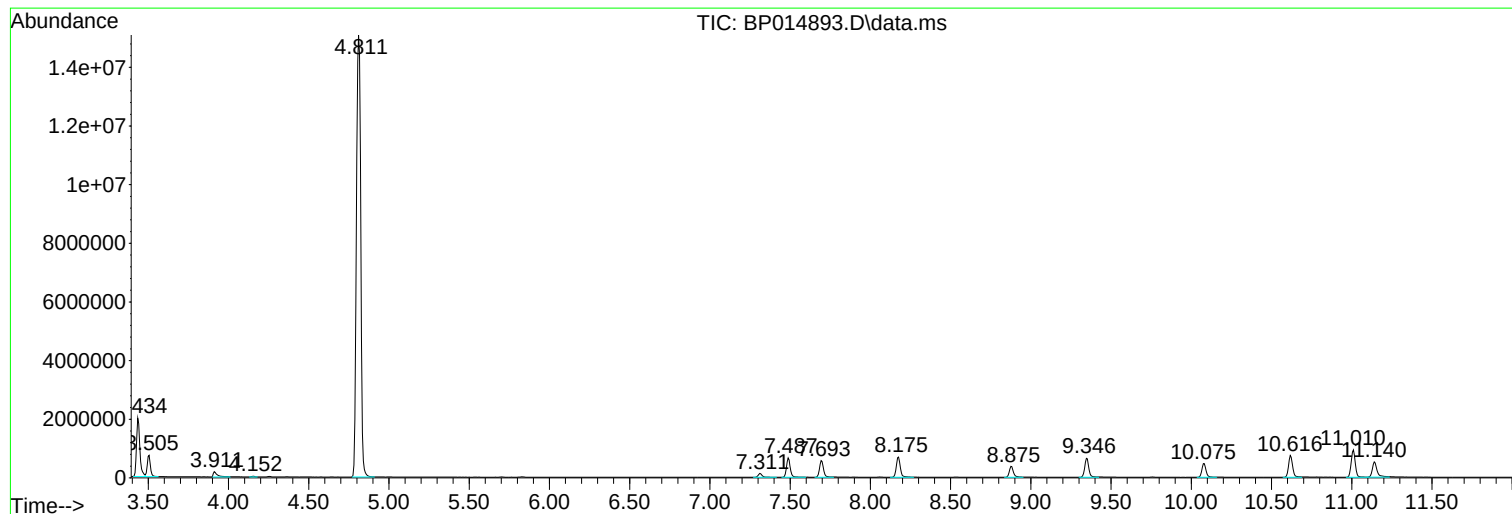
Sum of corrected areas: 76480914

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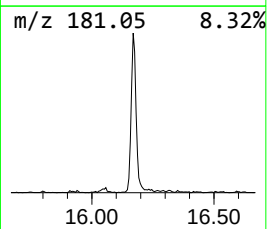
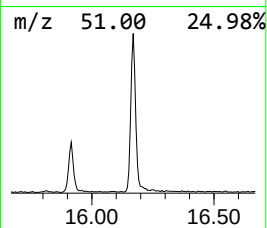
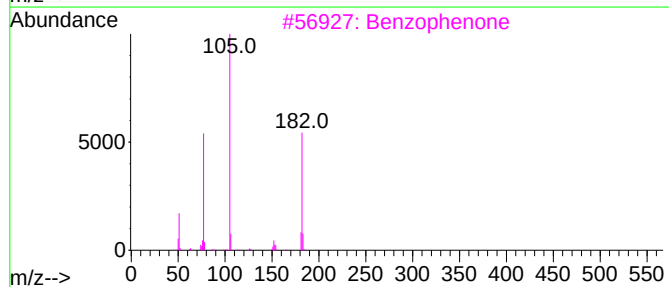
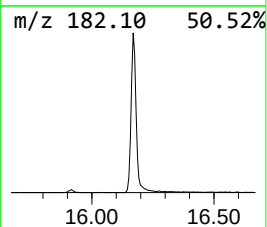
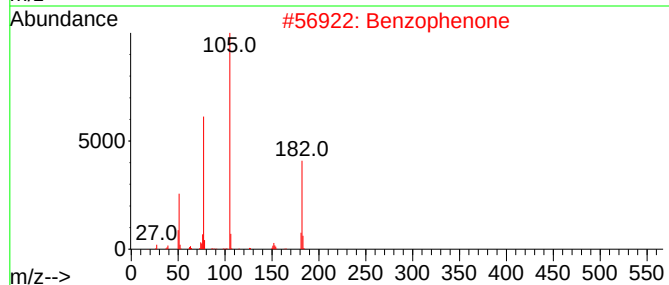
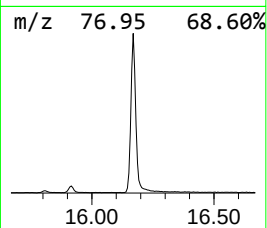
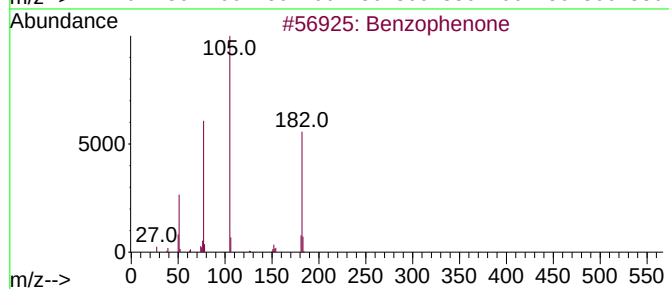
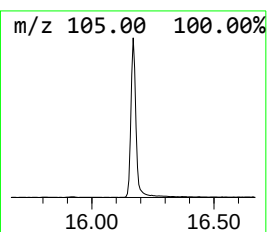
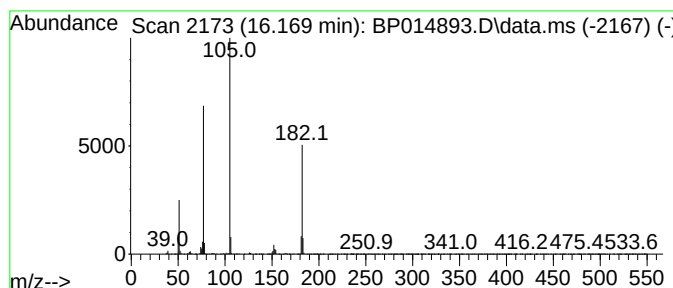
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042823.MA.M
Quant Title : SVOA CALIBRATION

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	7.73 ng/ul	789380	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	96
3			Benzophenone	182	C13H10O	000119-61-9	95
4			Benzophenone	182	C13H10O	000119-61-9	93
5			1,2,4,5-Tetroxane, 3,3,6,6-tetra...	396	C26H20O4	016204-36-7	78



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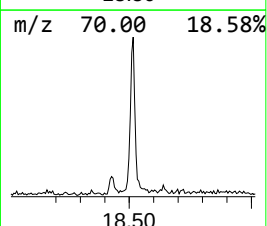
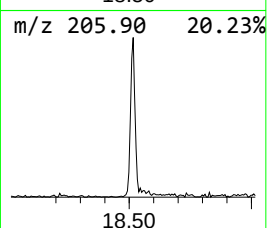
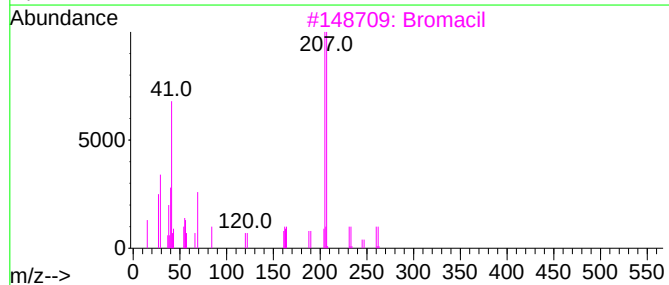
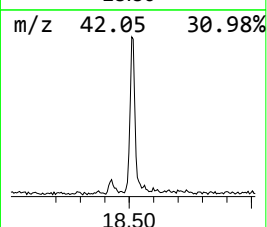
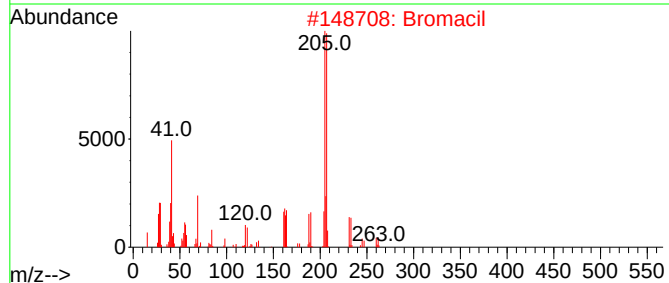
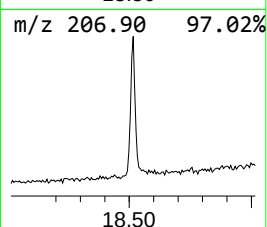
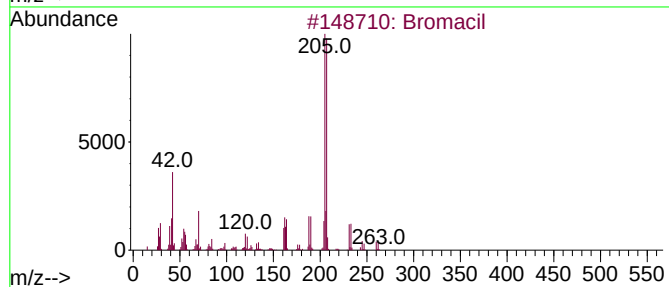
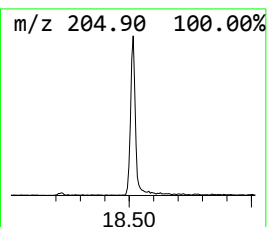
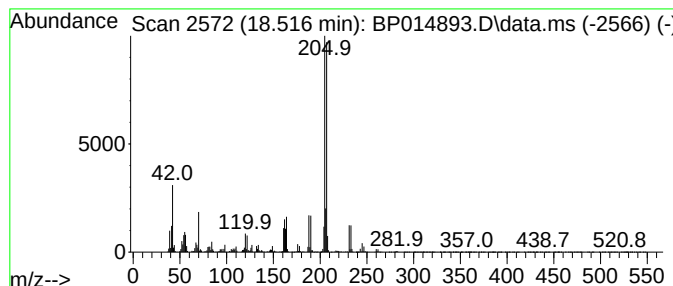
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Peak Number 3 Bromacil Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.516	3.05 ng/ul	335956	Phenanthrene-d10	17.575

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	78
4			5-Acetamido-2,4-dichlorobenzoic ...	247	C9H7Cl2NO3	050602-49-8	59
5			Benzoic acid, 3-amino-2,5-dichloro-	205	C7H5Cl2NO2	000133-90-4	50



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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	16.169	7.7	ng/ul	789380	3	14.816	2042250	20.0
Bromacil	18.516	3.0	ng/ul	335956	4	17.575	2201570	20.0