

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111423\
 Data File : BP018170.D
 Acq On : 14 Nov 2023 23:29
 Operator : MA/JU
 Sample : 05337-15
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GCDK0

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-BP111023.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP018170.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.164	10	13	16	rBV	14831	15889	0.59%	0.068%
2	3.617	87	90	104	rBV	26194	65376	2.45%	0.282%
3	3.781	113	118	123	rVB4	5313	10684	0.40%	0.046%
4	4.416	223	226	232	rBV8	3171	5241	0.20%	0.023%
5	5.828	462	466	468	rBV5	2500	3960	0.15%	0.017%
6	6.969	652	660	675	rVV	40943	99720	3.73%	0.430%
7	7.140	682	689	699	rBV	257459	418164	15.66%	1.802%
8	7.310	711	718	736	rBV	234925	432554	16.19%	1.864%
9	7.787	793	799	812	rBV	302975	516452	19.33%	2.225%
10	8.516	918	923	941	rBV	148187	320531	12.00%	1.381%
11	8.993	996	1004	1021	rBV	316727	591273	22.14%	2.548%
12	9.122	1023	1026	1027	rVV3	4684	5585	0.21%	0.024%
13	9.163	1030	1033	1037	rVB5	5215	8411	0.31%	0.036%
14	9.704	1118	1125	1146	rBV	263496	517648	19.38%	2.231%
15	9.851	1148	1150	1157	rVB8	1685	2766	0.10%	0.012%
16	10.240	1206	1216	1235	rBV	273226	644217	24.12%	2.776%
17	10.604	1271	1278	1286	rBV	424717	710325	26.59%	3.061%
18	10.793	1301	1310	1327	rBV	193532	480800	18.00%	2.072%
19	11.204	1376	1380	1385	rVB8	2357	4737	0.18%	0.020%
20	11.663	1453	1458	1462	rVB7	3733	5959	0.22%	0.026%
21	12.228	1551	1554	1555	rBV3	3678	3747	0.14%	0.016%
22	13.304	1731	1737	1745	rBV2	19025	33086	1.24%	0.143%
23	13.410	1751	1755	1756	rBV4	3428	3487	0.13%	0.015%
24	13.810	1820	1823	1827	rVB5	2241	2843	0.11%	0.012%
25	13.887	1830	1836	1858	rVV	882011	1295206	48.49%	5.581%
26	14.163	1875	1883	1894	rBV	849469	1247734	46.71%	5.377%
27	14.469	1928	1935	1944	rBV	630776	940204	35.20%	4.051%
28	14.798	1983	1991	1995	rBV10	13587	32105	1.20%	0.138%
29	14.875	2001	2004	2009	rVB7	3842	5987	0.22%	0.026%
30	15.204	2057	2060	2066	rVB4	4883	8343	0.31%	0.036%
31	15.251	2066	2068	2072	rVV5	2953	3929	0.15%	0.017%
32	15.298	2072	2076	2080	rVV7	2625	4874	0.18%	0.021%
33	15.334	2080	2082	2086	rVB5	2819	3152	0.12%	0.014%
34	15.469	2096	2105	2114	rBV	1184452	1695695	63.48%	7.307%
35	15.534	2114	2116	2121	rVB6	8291	11884	0.44%	0.051%
36	15.628	2126	2132	2144	rBV	367572	614130	22.99%	2.646%

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Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
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37	15.886	2172	2176	2180	rBV2	7807	10349	0.39%	0.045%
38	15.922	2180	2182	2185	rVV4	3528	4376	0.16%	0.019%
39	15.975	2189	2191	2195	rVB5	3270	3773	0.14%	0.016%
40	16.022	2195	2199	2201	rVB5	2821	3711	0.14%	0.016%
41	16.051	2201	2204	2206	rBV4	2208	3005	0.11%	0.013%
42	16.116	2213	2215	2222	rVB7	2241	3603	0.13%	0.016%
43	16.192	2222	2228	2233	rBV	22984	34764	1.30%	0.150%
44	16.281	2237	2243	2247	rBV	60455	86205	3.23%	0.371%
45	16.339	2248	2253	2259	rVV3	76841	150568	5.64%	0.649%
46	16.398	2259	2263	2268	rVV2	61972	112976	4.23%	0.487%
47	16.445	2268	2271	2276	rVV2	42658	60525	2.27%	0.261%
48	16.504	2276	2281	2283	rVV2	24126	43157	1.62%	0.186%
49	16.551	2286	2289	2293	rVV	24460	33461	1.25%	0.144%
50	16.604	2293	2298	2305	rVV4	53943	110618	4.14%	0.477%
51	16.675	2305	2310	2315	rVV	58921	94077	3.52%	0.405%
52	16.716	2315	2317	2319	rVV3	17460	20428	0.76%	0.088%
53	16.745	2319	2322	2329	rVB	29923	44297	1.66%	0.191%
54	16.886	2341	2346	2349	rBV6	3380	5153	0.19%	0.022%
55	16.916	2349	2351	2356	rVB6	3699	4320	0.16%	0.019%
56	16.998	2362	2365	2369	rVB5	3514	5388	0.20%	0.023%
57	17.051	2371	2374	2376	rVB4	2624	2781	0.10%	0.012%
58	17.086	2376	2380	2383	rBV6	3348	4735	0.18%	0.020%
59	17.122	2384	2386	2389	rVB3	3935	3774	0.14%	0.016%
60	17.198	2394	2399	2401	rBV6	2564	4939	0.18%	0.021%
61	17.245	2401	2407	2418	rBV	847553	1183693	44.32%	5.101%
62	17.351	2418	2425	2439	rBV	1396640	2065133	77.31%	8.899%
63	17.510	2447	2452	2455	rBV5	5540	10643	0.40%	0.046%
64	17.580	2461	2464	2468	rVB6	3051	4908	0.18%	0.021%
65	17.880	2510	2515	2522	rBV2	57322	76368	2.86%	0.329%
66	18.098	2548	2552	2561	rBV2	38281	62752	2.35%	0.270%
67	18.192	2564	2568	2572	rVB	9010	11861	0.44%	0.051%
68	18.310	2584	2588	2600	rBV	24651	40255	1.51%	0.173%
69	18.680	2649	2651	2654	rVB4	3372	3330	0.12%	0.014%
70	19.180	2731	2736	2739	rBV	19708	27611	1.03%	0.119%
71	19.233	2741	2745	2753	rVV5	29673	71257	2.67%	0.307%
72	19.345	2761	2764	2770	rVB3	21476	27929	1.05%	0.120%
73	19.480	2784	2787	2790	rVB4	10373	10247	0.38%	0.044%
74	19.604	2802	2808	2817	rBV	2069832	2600043	97.34%	11.204%
75	20.510	2960	2962	2966	rVB5	11556	10386	0.39%	0.045%
76	21.380	3104	3110	3116	rBV	946341	1341205	50.21%	5.779%

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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-BP111023.MA.M
Title : SVOA CALIBRATION

77	23.680	3494	3501	3515	rVB2	1314424	2671086	100.00%	11.510%
78	23.845	3523	3529	3542	rVB	659186	1370207	51.30%	5.904%

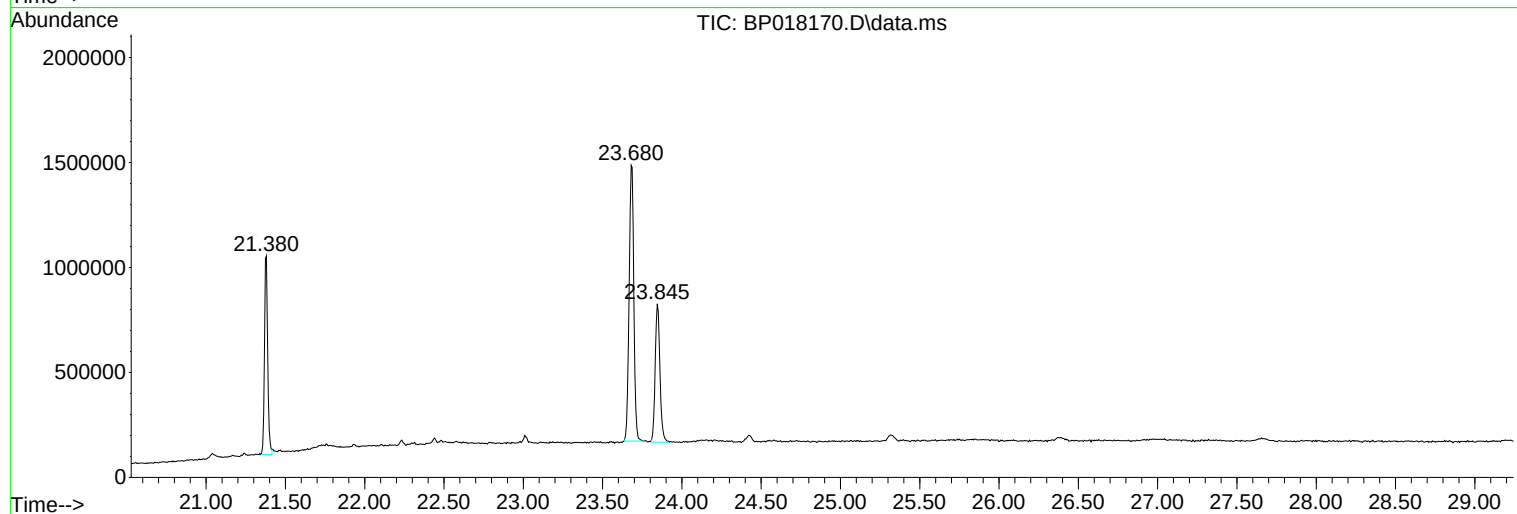
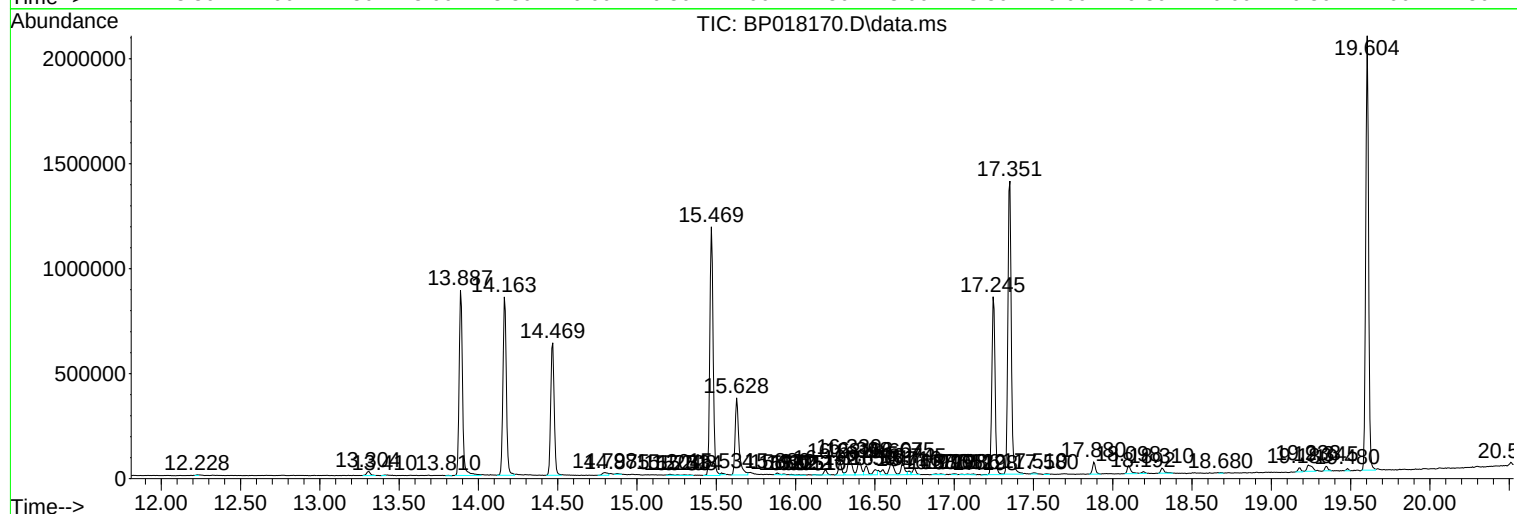
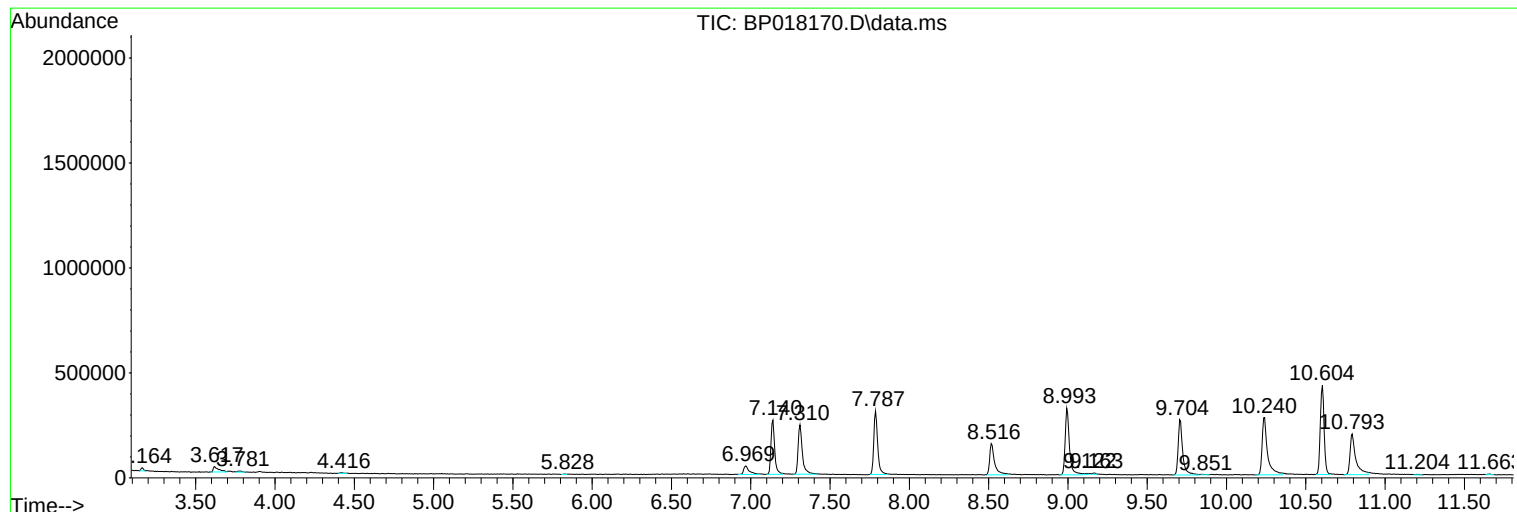
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

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



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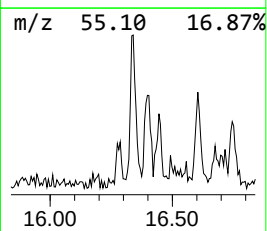
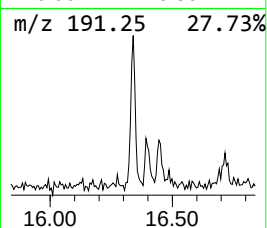
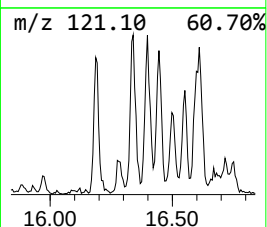
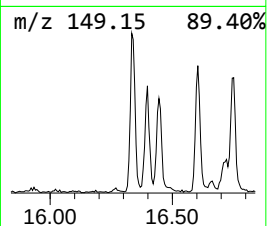
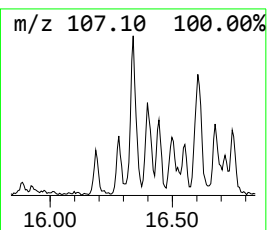
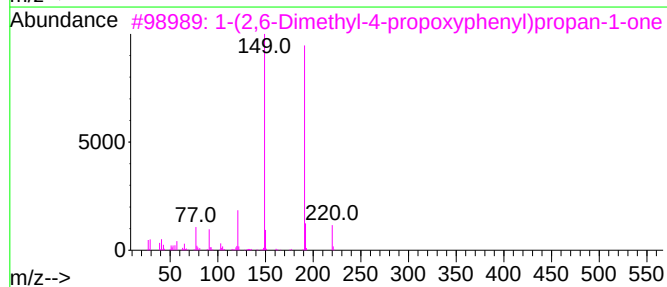
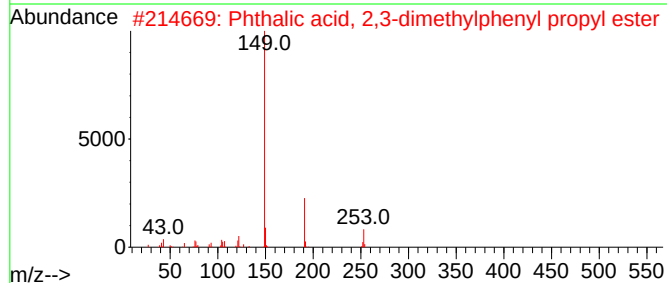
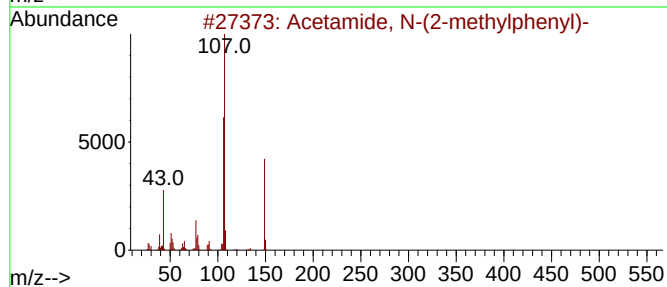
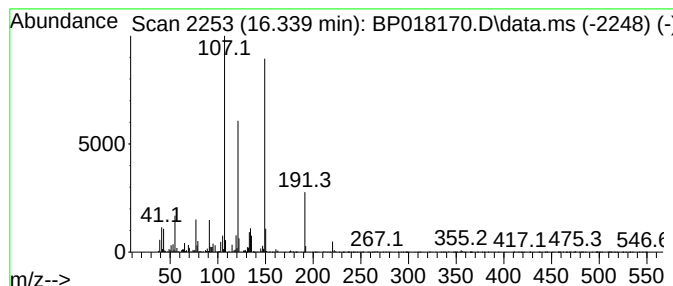
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 unknown-01 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.339	2.54 ng/ul	150568	Phenanthrene-d10	17.245

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acetamide, N-(2-methylphenyl)-	149	C9H11NO	000120-66-1	46
2			Phthalic acid, 2,3-dimethylpheny...	312	C19H20O4	1000357-08-5	38
3			1-(2,6-Dimethyl-4-propoxyphenyl)...	220	C14H20O2	1000190-22-3	38
4			Acetamide, N-(4-methylphenyl)-	149	C9H11NO	000103-89-9	38
5			Methyl (1s,3s,5R,7S)-3-methylada...	208	C13H20O2	1010504-39-1	35



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TIC Library : C:\DATABASE\NIST20.L
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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown-01	16.339	2.5	ng/u1	150568	4	17.245	1183690	20.0