

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110924\
 Data File : BP022984.D
 Acq On : 10 Nov 2024 03:26
 Operator : RC/JU
 Sample : P4746-15
 Misc :
 ALS Vial : 71 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GCP88

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-BP110924.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP022984.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.111	16	21	24	rBV	54376	75988	2.63%	0.340%
2	3.140	24	26	32	rVB	14608	17739	0.61%	0.079%
3	3.564	94	98	114	rBV	40945	79842	2.76%	0.358%
4	3.864	146	149	159	rVB4	3935	6400	0.22%	0.029%
5	4.234	208	212	217	rVB8	2189	2932	0.10%	0.013%
6	4.375	229	236	244	rVB4	11370	22288	0.77%	0.100%
7	4.934	329	331	339	rBV7	2025	3429	0.12%	0.015%
8	6.281	555	560	566	rBV8	1560	2976	0.10%	0.013%
9	6.793	641	647	660	rBV2	59513	128370	4.44%	0.575%
10	6.934	665	671	680	rBV	193282	299331	10.34%	1.341%
11	7.134	698	705	715	rBV	211664	379999	13.13%	1.702%
12	7.205	715	717	726	rVB10	2868	5475	0.19%	0.025%
13	7.587	774	782	797	rBV	220851	383128	13.24%	1.716%
14	8.305	898	904	918	rBV	211613	370976	12.82%	1.661%
15	8.728	968	976	989	rBV	278777	459324	15.87%	2.057%
16	8.875	998	1001	1008	rVB4	2848	4159	0.14%	0.019%
17	9.122	1037	1043	1048	rVB5	2945	5434	0.19%	0.024%
18	9.446	1091	1098	1110	rBV	236441	414105	14.31%	1.855%
19	9.987	1181	1190	1211	rBV	291165	632436	21.86%	2.832%
20	10.163	1215	1220	1232	rVB8	4668	14228	0.49%	0.064%
21	10.340	1243	1250	1261	rBV	310526	527084	18.21%	2.361%
22	10.493	1267	1276	1290	rBV	202398	423838	14.65%	1.898%
23	11.598	1458	1464	1467	rBV7	1748	3044	0.11%	0.014%
24	11.946	1514	1523	1530	rBV2	4621	10684	0.37%	0.048%
25	13.616	1801	1807	1817	rBV	881368	1134930	39.22%	5.083%
26	13.904	1845	1856	1868	rBV2	818372	1148708	39.70%	5.144%
27	14.216	1902	1909	1917	rBV	547027	838844	28.99%	3.757%
28	14.510	1946	1959	1972	rBV4	36243	133194	4.60%	0.597%
29	15.222	2071	2080	2092	rBV	863593	1605359	55.48%	7.190%
30	15.369	2099	2105	2116	rBV	397537	599283	20.71%	2.684%
31	15.457	2118	2120	2128	rVB5	6478	9753	0.34%	0.044%
32	15.598	2139	2144	2151	rBV	27318	39796	1.38%	0.178%
33	16.486	2292	2295	2300	rVB7	2324	3594	0.12%	0.016%
34	16.604	2311	2315	2321	rVB2	4829	6422	0.22%	0.029%
35	16.686	2326	2329	2336	rVB9	1605	2942	0.10%	0.013%
36	16.969	2369	2377	2378	rBV5	5199	7632	0.26%	0.034%

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37	17.016	2378	2385	2396	rVV	753509	1192018	41.19%	5.338%
38	17.122	2397	2403	2417	rVV	1420927	1973678	68.21%	8.839%
39	17.445	2450	2458	2463	rBV	2490	6237	0.22%	0.028%
40	17.516	2464	2470	2480	rBV	244130	373850	12.92%	1.674%
41	17.675	2494	2497	2500	rBV4	5503	7724	0.27%	0.035%
42	17.710	2500	2503	2511	rVB8	9772	15073	0.52%	0.068%
43	17.886	2529	2533	2537	rVB7	2136	3113	0.11%	0.014%
44	17.951	2539	2544	2550	rBV	33308	48304	1.67%	0.216%
45	18.163	2574	2580	2590	rBV	128866	209106	7.23%	0.936%
46	19.033	2724	2728	2735	rVB4	23703	34477	1.19%	0.154%
47	19.122	2738	2743	2748	rBV	16254	26907	0.93%	0.121%
48	19.292	2768	2772	2777	rBV8	12895	20337	0.70%	0.091%
49	19.439	2794	2797	2799	rBV4	4471	6061	0.21%	0.027%
50	19.498	2800	2807	2816	rVV	1822810	2564520	88.62%	11.485%
51	21.445	3132	3138	3149	rBV	994376	1546214	53.43%	6.925%
52	24.445	3637	3648	3659	rVB	1114978	2893717	100.00%	12.960%
53	24.651	3676	3683	3697	rVB	650758	1603893	55.43%	7.183%

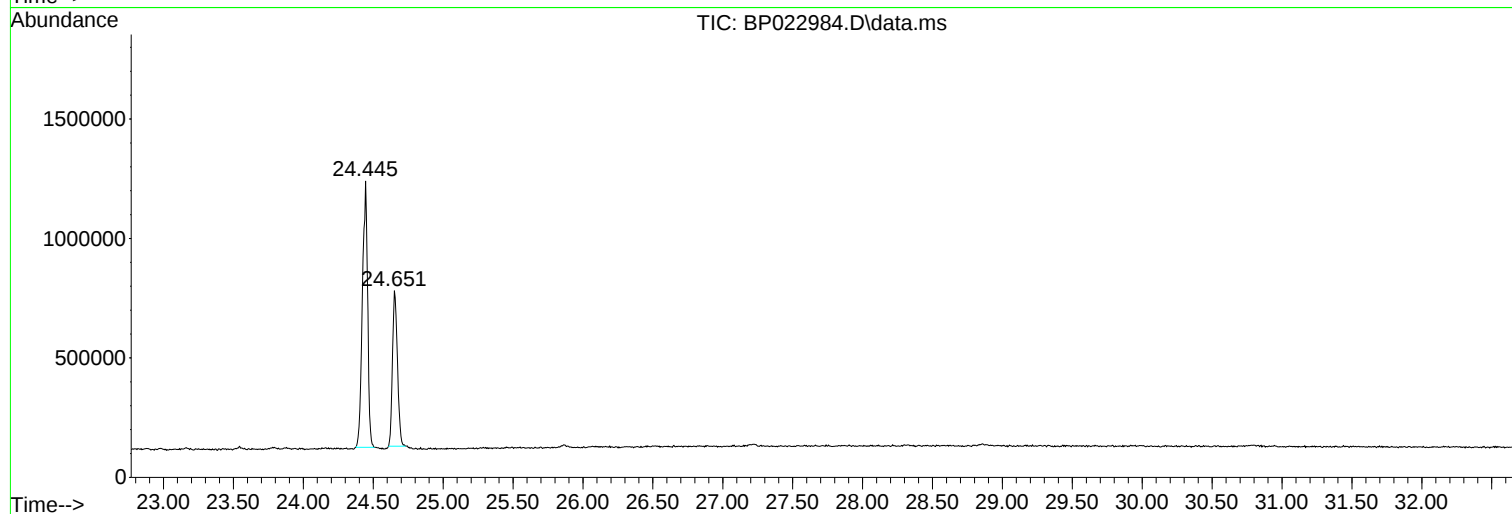
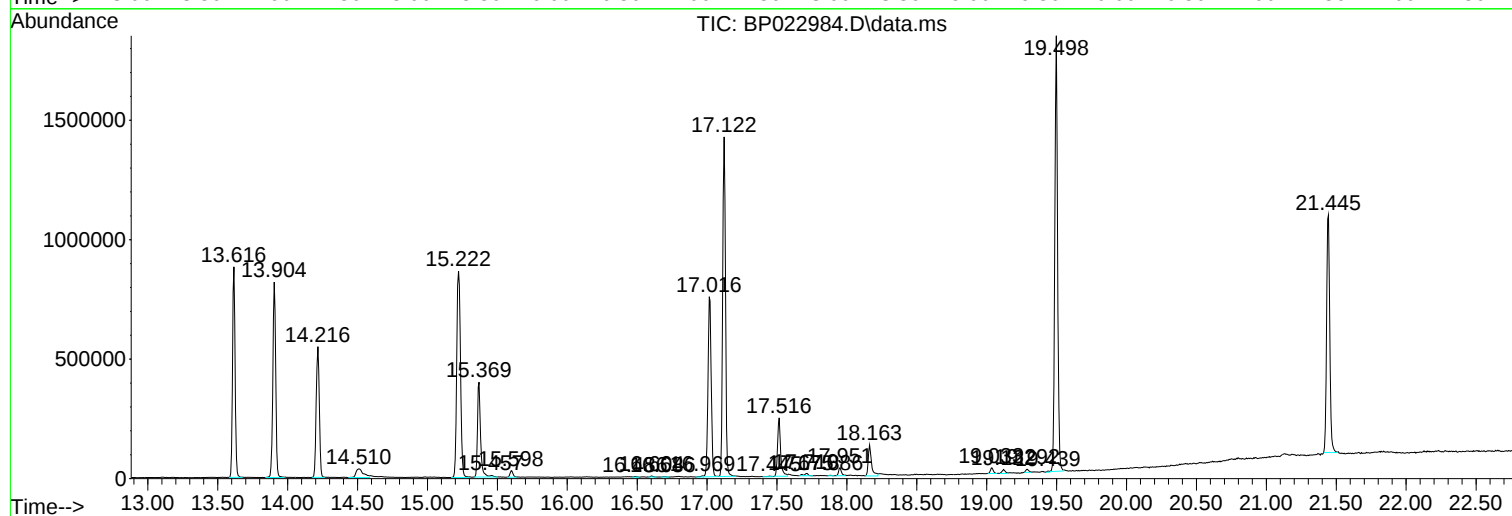
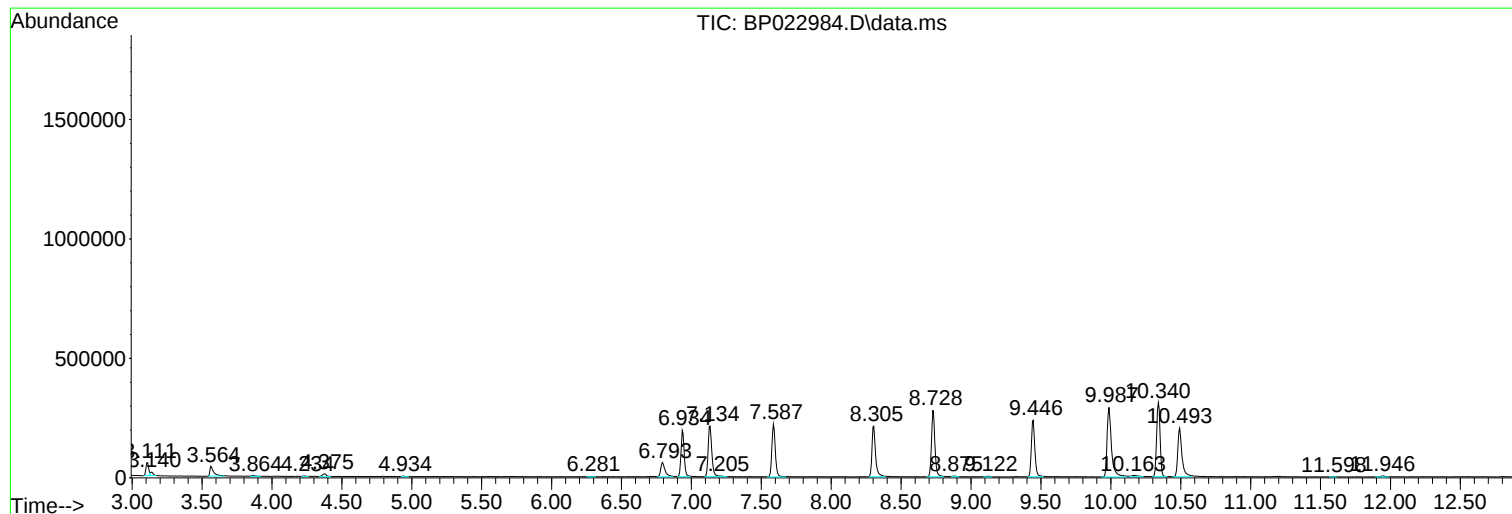
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110924.MA.M
Quant Title : SVOA CALIBRATION

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



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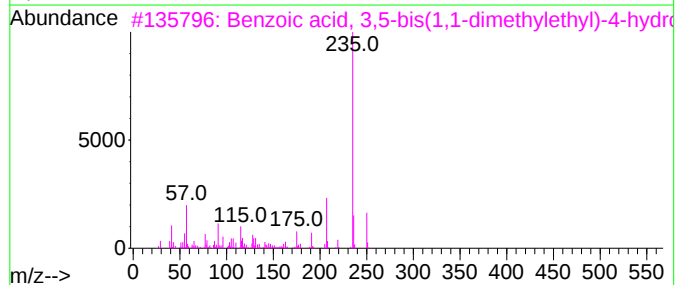
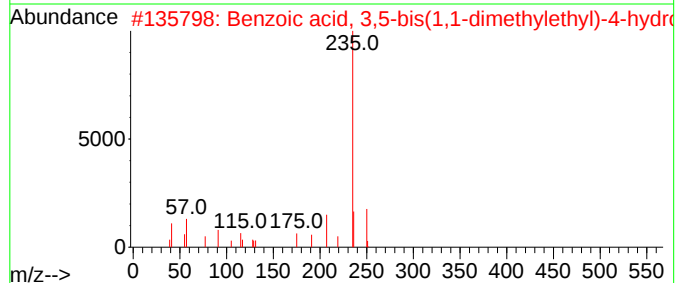
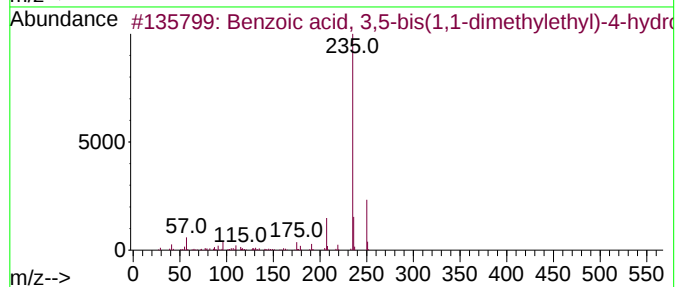
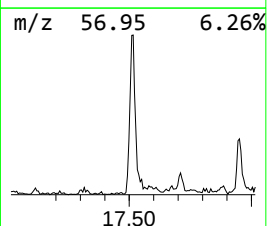
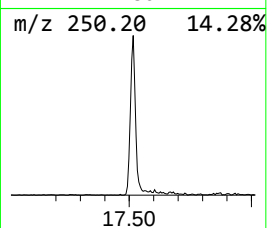
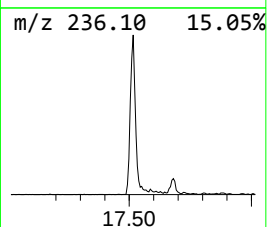
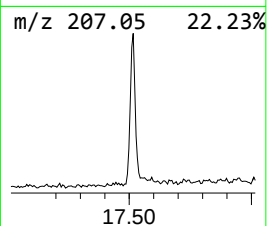
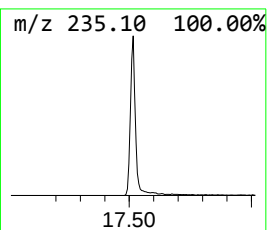
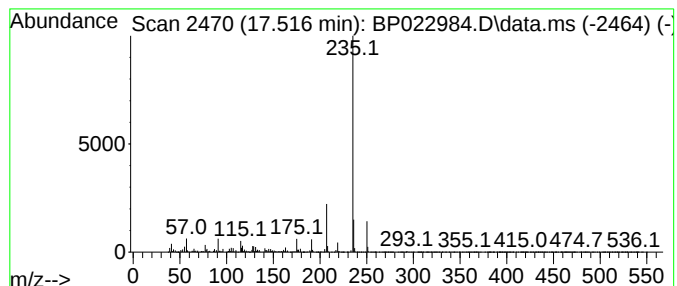
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzoic acid, 3,5-bis(1,1-d... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.516	6.27 ng/ul	373850	Phenanthrene-d10	17.016

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzoic acid, 3,5-bis(1,1-dimeth...	250	C15H22O3	001421-49-4	95
2			Benzoic acid, 3,5-bis(1,1-dimeth...	250	C15H22O3	001421-49-4	94
3			Benzoic acid, 3,5-bis(1,1-dimeth...	250	C15H22O3	001421-49-4	94
4			Ethane, 1,1-bis(p-chlorophenyl)-	250	C14H12Cl2	003547-04-4	53
5			N-(4-Hydroxyphenyl)-2-naphthylamine	235	C16H13NO	000093-45-8	53



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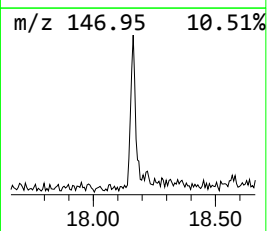
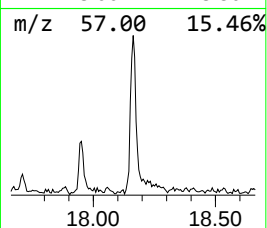
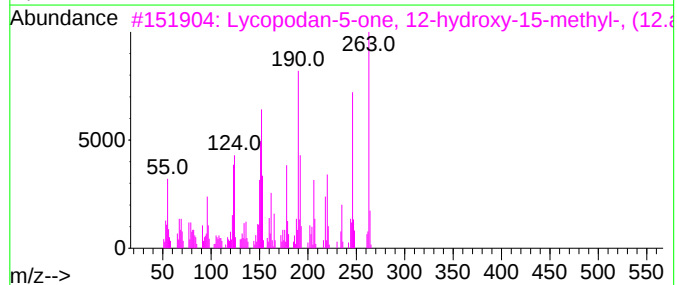
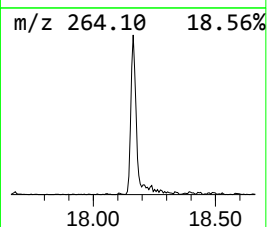
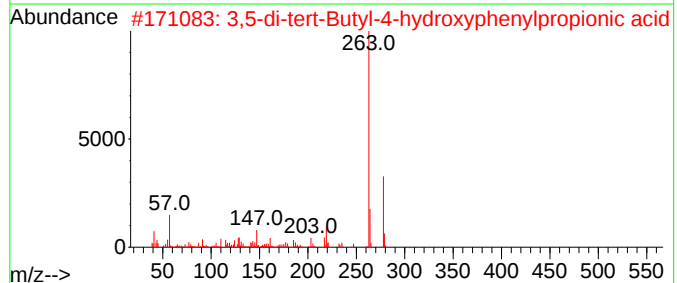
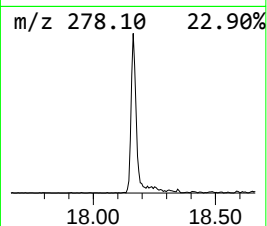
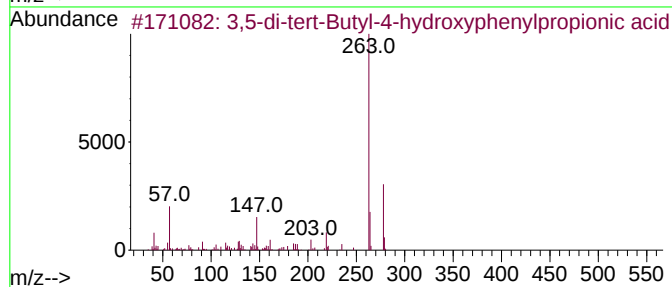
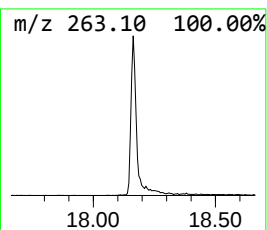
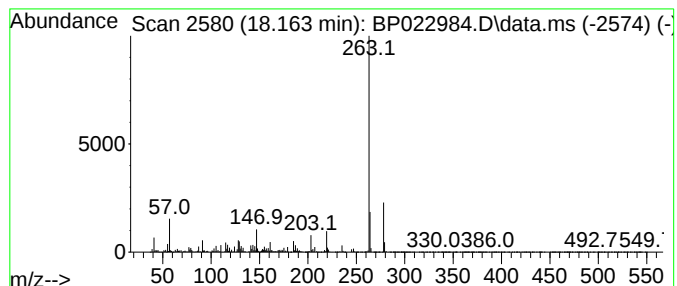
Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 3,5-di-tert-Butyl-4-hydroxy... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.163	3.51 ng/ul	209106	Phenanthrene-d10	17.016	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3,5-di-tert-Butyl-4-hydroxypheny...	278	C17H26O3	020170-32-5	94
2	3,5-di-tert-Butyl-4-hydroxypheny...	278	C17H26O3	020170-32-5	94
3	Lycopodan-5-one, 12-hydroxy-15-m...	263	C16H25NO2	021061-90-5	91
4	6,7-Dimethoxy-4(1H)-quinazolinon...	278	C13H18N2O3Si	1000479-65-5	64
5	2,4-Di-tert-butylphenoxytrimethy...	278	C17H30O5Si	053925-65-8	64



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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
Benzoic acid, 3...	17.516	6.3	ng/ul	373850	4	17.016	1192020	20.0
3,5-di-tert-But...	18.163	3.5	ng/ul	209106	4	17.016	1192020	20.0