

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110924\
 Data File : BP022974.D
 Acq On : 09 Nov 2024 19:57
 Operator : RC/JU
 Sample : P4746-19
 Misc :
 ALS Vial : 59 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GCP92

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: BP022974.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.146	23	27	31	rVB	10960	13416	0.52%	0.070%
2	3.569	95	99	115	rBV	30285	59615	2.31%	0.311%
3	3.869	146	150	155	rVB4	2860	4489	0.17%	0.023%
4	6.805	642	649	664	rBV	58508	123552	4.78%	0.644%
5	6.940	666	672	683	rBV	158819	255614	9.89%	1.333%
6	7.140	699	706	716	rBV	198473	343824	13.30%	1.793%
7	7.593	776	783	793	rBV	209921	337637	13.06%	1.760%
8	8.299	897	903	918	rBV	186341	337914	13.07%	1.762%
9	8.734	968	977	997	rBV	219993	388010	15.01%	2.023%
10	9.134	1036	1045	1048	rBV9	1956	3329	0.13%	0.017%
11	9.322	1067	1077	1082	rBV10	1598	3715	0.14%	0.019%
12	9.446	1091	1098	1110	rBV	182973	345254	13.36%	1.800%
13	9.993	1180	1191	1211	rBV	244048	522512	20.21%	2.724%
14	10.340	1241	1250	1268	rVB	281151	466926	18.06%	2.434%
15	10.493	1268	1276	1294	rBV2	198409	434593	16.81%	2.266%
16	11.957	1520	1525	1534	rBV5	2903	6577	0.25%	0.034%
17	13.098	1715	1719	1726	rBV7	2510	4096	0.16%	0.021%
18	13.610	1800	1806	1821	rBV	693588	917857	35.50%	4.785%
19	13.739	1821	1828	1835	rVB10	3991	8796	0.34%	0.046%
20	13.898	1849	1855	1868	rVB	690966	915994	35.43%	4.776%
21	14.210	1900	1908	1919	rBV	479537	712203	27.55%	3.713%
22	14.504	1952	1958	1962	rBV2	33087	66868	2.59%	0.349%
23	14.681	1986	1988	1994	rVB7	2657	3499	0.14%	0.018%
24	15.228	2072	2081	2093	rBV	758359	1325457	51.27%	6.910%
25	15.369	2099	2105	2116	rBV	341673	466514	18.05%	2.432%
26	15.451	2117	2119	2131	rVB9	5473	12465	0.48%	0.065%
27	15.598	2139	2144	2153	rVB	66262	98915	3.83%	0.516%
28	16.163	2236	2240	2246	rVB7	4526	7257	0.28%	0.038%
29	16.616	2312	2317	2323	rBV9	2044	4178	0.16%	0.022%
30	17.016	2375	2385	2394	rBV	639320	973533	37.66%	5.076%
31	17.116	2396	2402	2414	rVV	1154671	1619475	62.64%	8.443%
32	17.504	2463	2468	2477	rBV	66598	94811	3.67%	0.494%
33	17.710	2499	2503	2511	rVB7	11017	17029	0.66%	0.089%
34	17.822	2518	2522	2523	rBV3	2321	2757	0.11%	0.014%
35	17.945	2539	2543	2552	rBV	34355	52272	2.02%	0.273%
36	18.169	2574	2581	2601	rBV	394236	648226	25.07%	3.380%

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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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37	19.033	2722	2728	2736	rBV7	18857	37791	1.46%	0.197%
38	19.122	2738	2743	2747	rBV3	11018	19205	0.74%	0.100%
39	19.292	2769	2772	2780	rVB9	10818	17140	0.66%	0.089%
40	19.492	2799	2806	2816	rBV	1412950	2088781	80.80%	10.890%
41	21.451	3133	3139	3147	rBV	849481	1336947	51.72%	6.970%
42	24.451	3640	3649	3663	rVB	958325	2585169	100.00%	13.478%
43	24.668	3677	3686	3697	rBV	530628	1496451	57.89%	7.802%

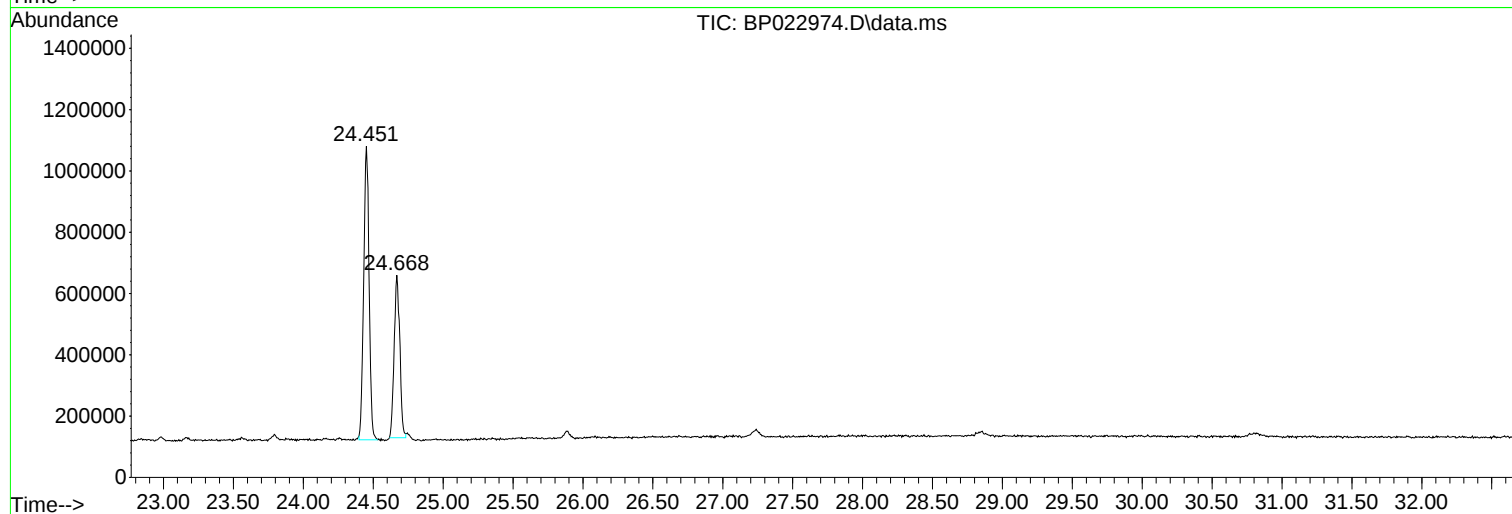
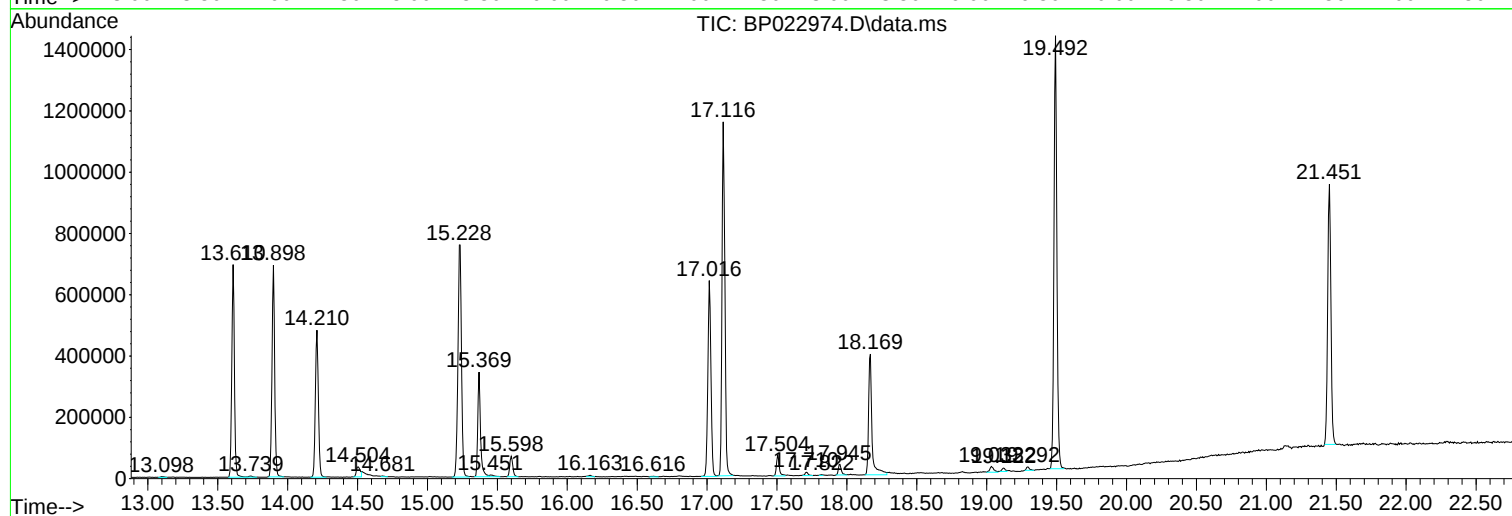
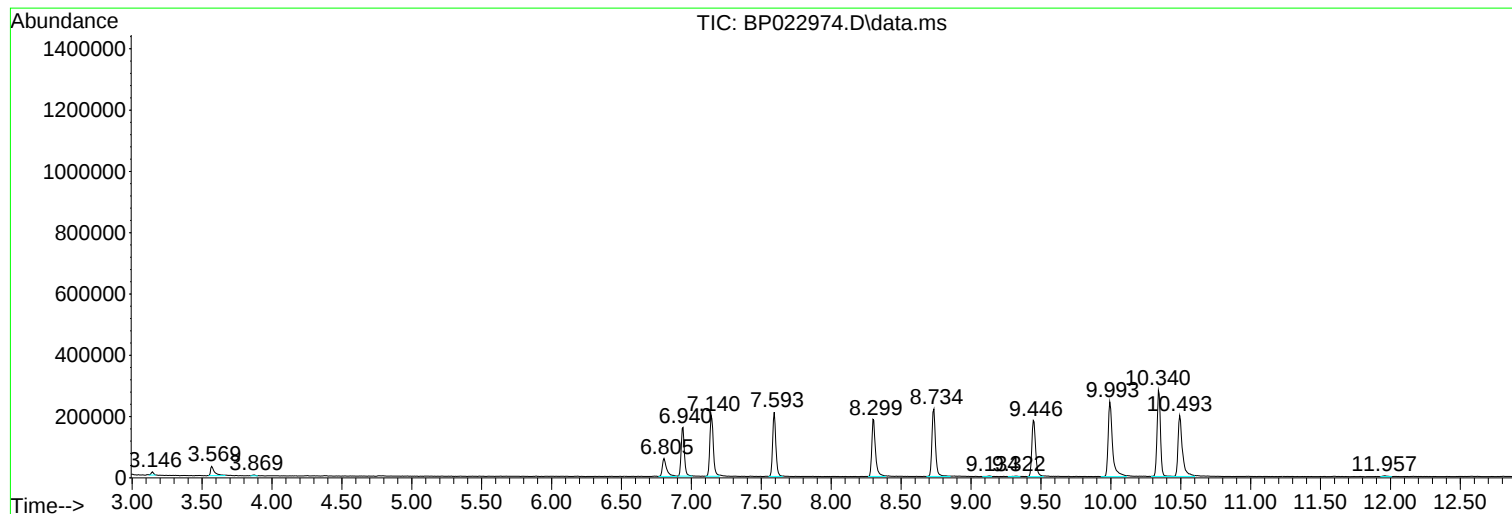
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110924.MA.M
Quant Title : SVOA CALIBRATION

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



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Instrument :
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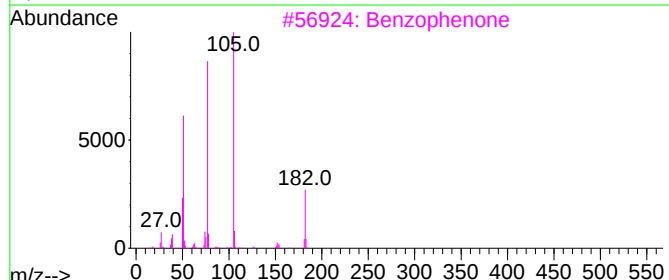
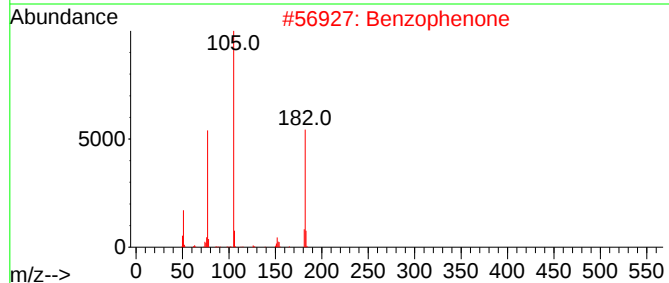
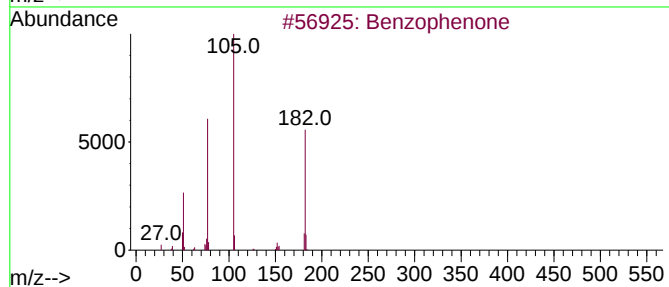
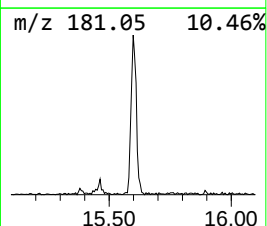
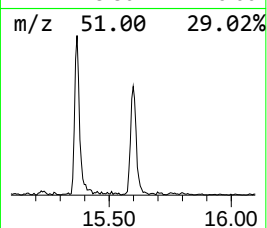
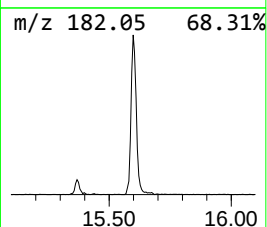
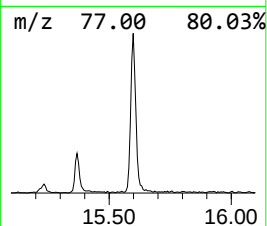
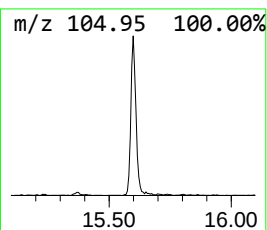
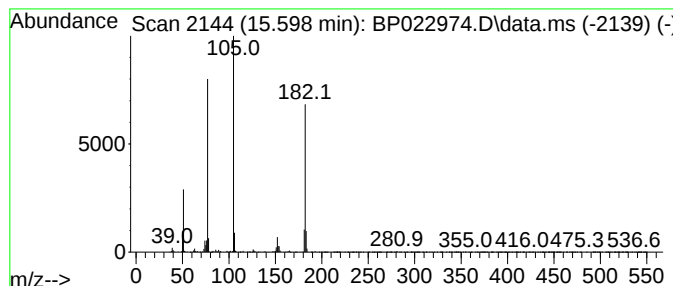
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110924.MA.M
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TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Benzophenone Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.598	2.78 ng/ul	98915	Acenaphthene-d10	14.210

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	89
4			Benzophenone	182	C13H10O	000119-61-9	87
5			1,2,4-Trioxolane, 3,3,5-triphenyl-	304	C20H16O3	023246-12-0	74



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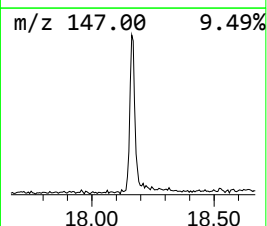
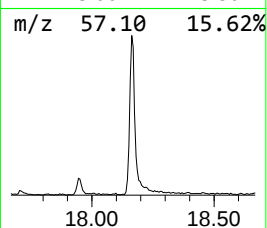
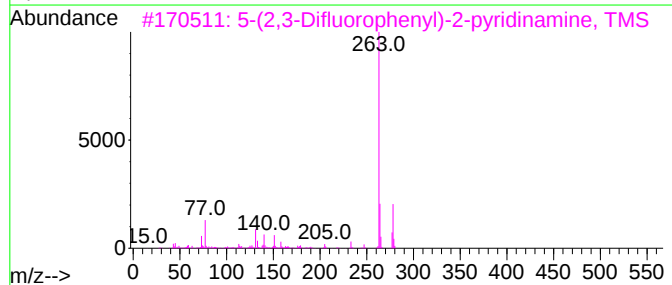
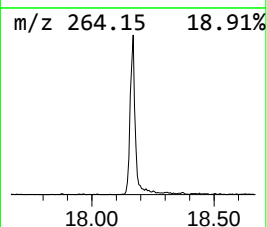
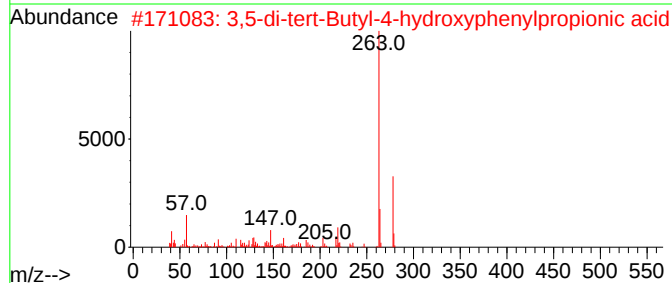
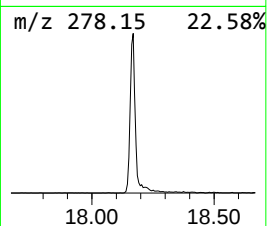
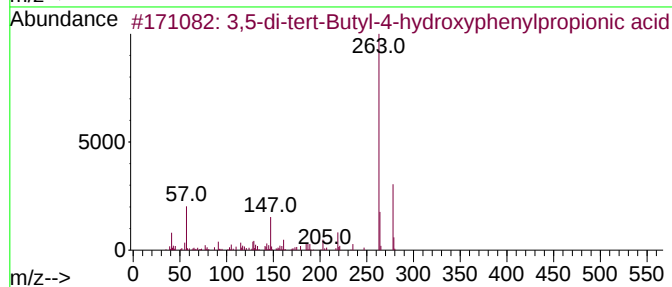
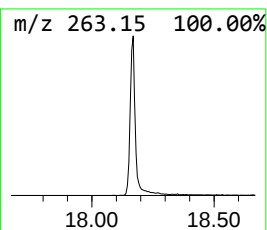
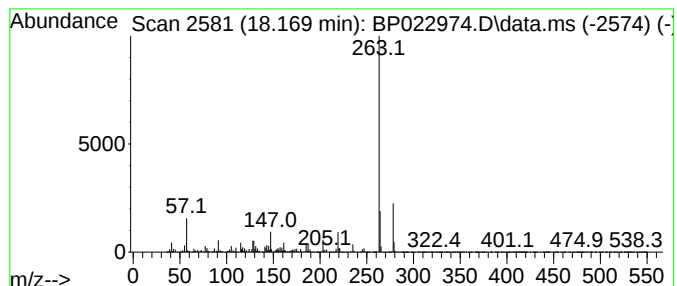
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TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 3,5-di-tert-Butyl-4-hydroxy... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD		R.T.		
18.169	13.32 ng/ul	648226	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	91
3			5-(2,3-Difluorophenyl)-2-pyridin...	278	C14H16F2N2Si	1000504-49-4	72
4			1H-Isoindole-1,3(2H)-dione, 2-(2...	263	C17H13NO2	016307-59-8	60
5			Lycopodan-5-one, 12-hydroxy-15-m...	263	C16H25NO2	021061-90-5	60



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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	15.598	2.8	ng/u1	98915	3	14.210	712203	20.0
3,5-di-tert-But...	18.169	13.3	ng/u1	648226	4	17.016	973533	20.0