

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110924\
 Data File : BP022988.D
 Acq On : 10 Nov 2024 06:10
 Operator : RC/JU
 Sample : P4746-08
 Misc :
 ALS Vial : 75 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GCP83

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: BP022988.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.140	22	26	33	rVB	15520	20939	0.64%	0.094%
2	3.558	94	97	116	rBV	50940	99001	3.01%	0.443%
3	3.858	145	148	155	rVB5	4025	6236	0.19%	0.028%
4	4.364	228	234	240	rBV2	6942	11432	0.35%	0.051%
5	6.734	629	637	641	rBV7	2305	4287	0.13%	0.019%
6	6.799	641	648	661	rVV	73803	153991	4.68%	0.689%
7	6.934	665	671	686	rVB	219008	348974	10.60%	1.561%
8	7.134	699	705	724	rBV	267117	452756	13.76%	2.025%
9	7.587	773	782	796	rBV	196259	332361	10.10%	1.487%
10	8.293	895	902	919	rBV	223544	443150	13.47%	1.982%
11	8.734	969	977	993	rBV	285337	523980	15.92%	2.344%
12	9.128	1038	1044	1049	rVB4	6205	9277	0.28%	0.042%
13	9.440	1090	1097	1112	rBV	246864	470078	14.28%	2.103%
14	9.981	1183	1189	1213	rBV	351146	712573	21.65%	3.188%
15	10.334	1242	1249	1261	rVB	270050	448656	13.63%	2.007%
16	10.487	1267	1275	1297	rBV	257172	562194	17.08%	2.515%
17	11.599	1456	1464	1473	rVB3	14736	24474	0.74%	0.109%
18	11.946	1517	1523	1530	rBV4	4349	9294	0.28%	0.042%
19	13.604	1797	1805	1820	rBV	693245	1211939	36.83%	5.422%
20	13.887	1846	1853	1866	rBV2	798800	1242782	37.77%	5.560%
21	14.204	1899	1907	1922	rBV	470828	707849	21.51%	3.167%
22	14.493	1949	1956	1975	rBV3	38465	147304	4.48%	0.659%
23	15.016	2042	2045	2051	rBV4	2706	4407	0.13%	0.020%
24	15.222	2073	2080	2096	rBV	1412299	1741180	52.91%	7.789%
25	15.363	2097	2104	2120	rBV	356869	630940	19.17%	2.823%
26	15.475	2120	2123	2129	rVB3	7285	10082	0.31%	0.045%
27	16.610	2312	2316	2319	rBV5	3347	4664	0.14%	0.021%
28	17.010	2378	2384	2393	rVV	782435	958475	29.13%	4.288%
29	17.104	2393	2400	2417	rVB2	1396829	2100339	63.82%	9.396%
30	17.310	2433	2435	2440	rVB4	2368	3545	0.11%	0.016%
31	17.704	2498	2502	2505	rBV6	3355	3928	0.12%	0.018%
32	17.934	2537	2541	2549	rBV4	22363	37876	1.15%	0.169%
33	18.010	2552	2554	2558	rVV3	3388	4536	0.14%	0.020%
34	18.157	2572	2579	2594	rBV	77218	147577	4.48%	0.660%
35	19.034	2724	2728	2735	rBV3	16303	29818	0.91%	0.133%
36	19.128	2739	2744	2747	rBV	20160	32754	1.00%	0.147%

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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

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37	19.286	2767	2771	2776	rBV5	27434	35031	1.06%	0.157%
38	19.498	2801	2807	2815	rBV	1586984	2652135	80.59%	11.865%
39	21.451	3132	3139	3147	rBV	823926	1302563	39.58%	5.827%
40	24.445	3639	3648	3665	rVB	1262071	3290822	100.00%	14.722%
41	24.674	3676	3687	3700	rVB	486541	1418897	43.12%	6.348%

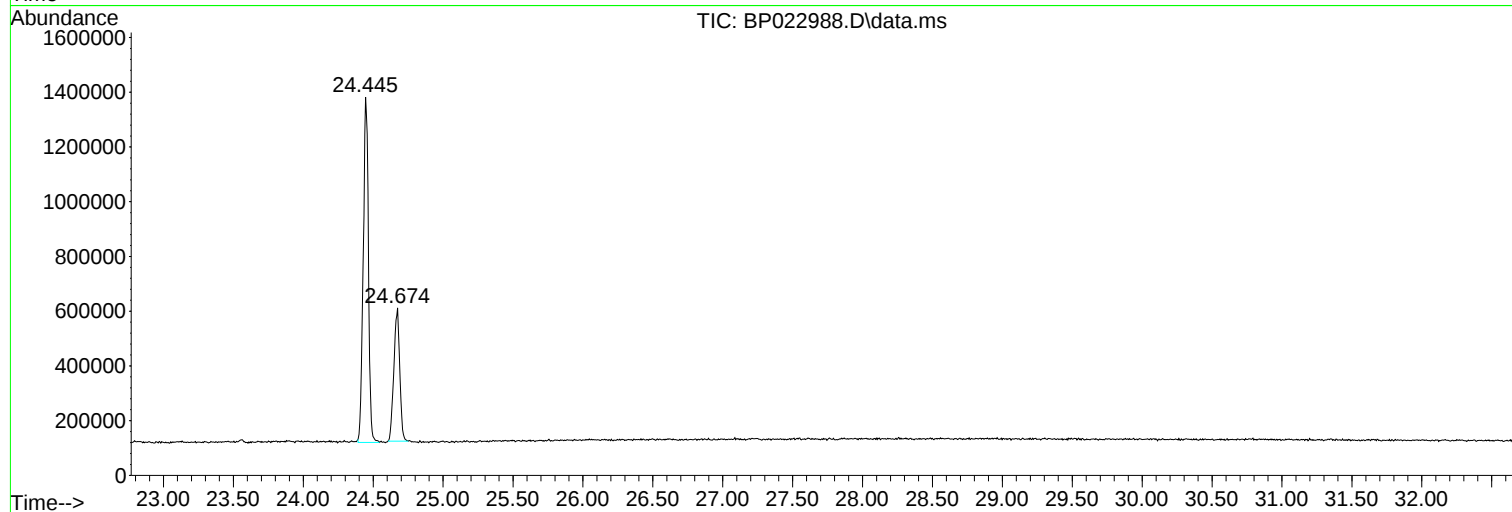
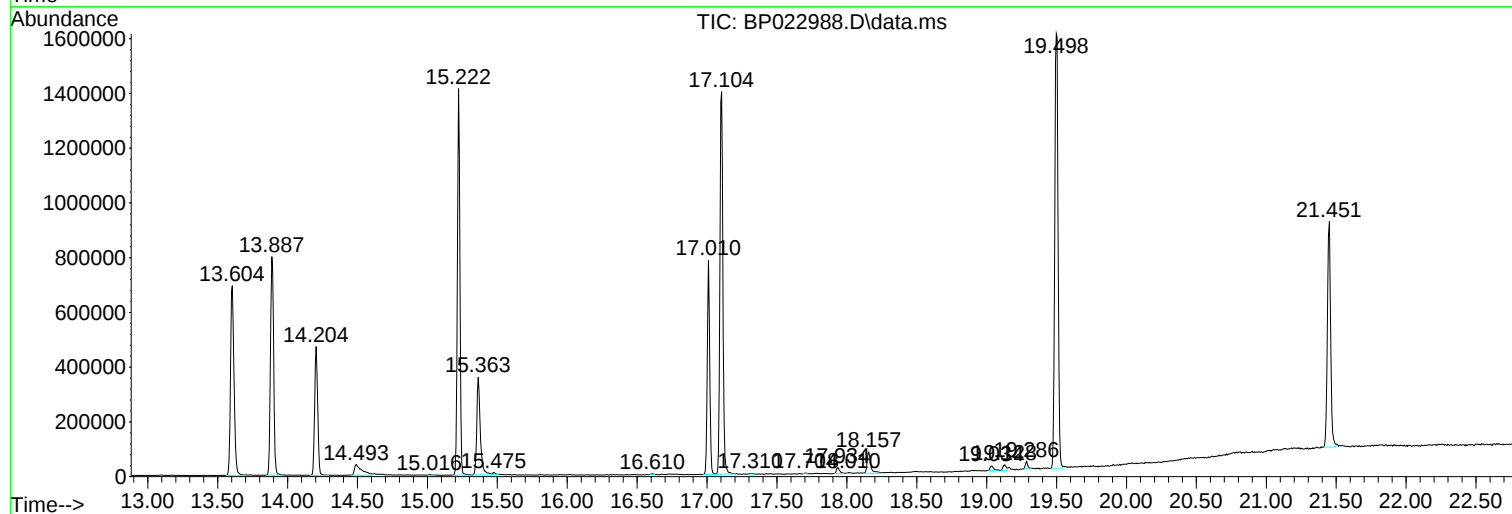
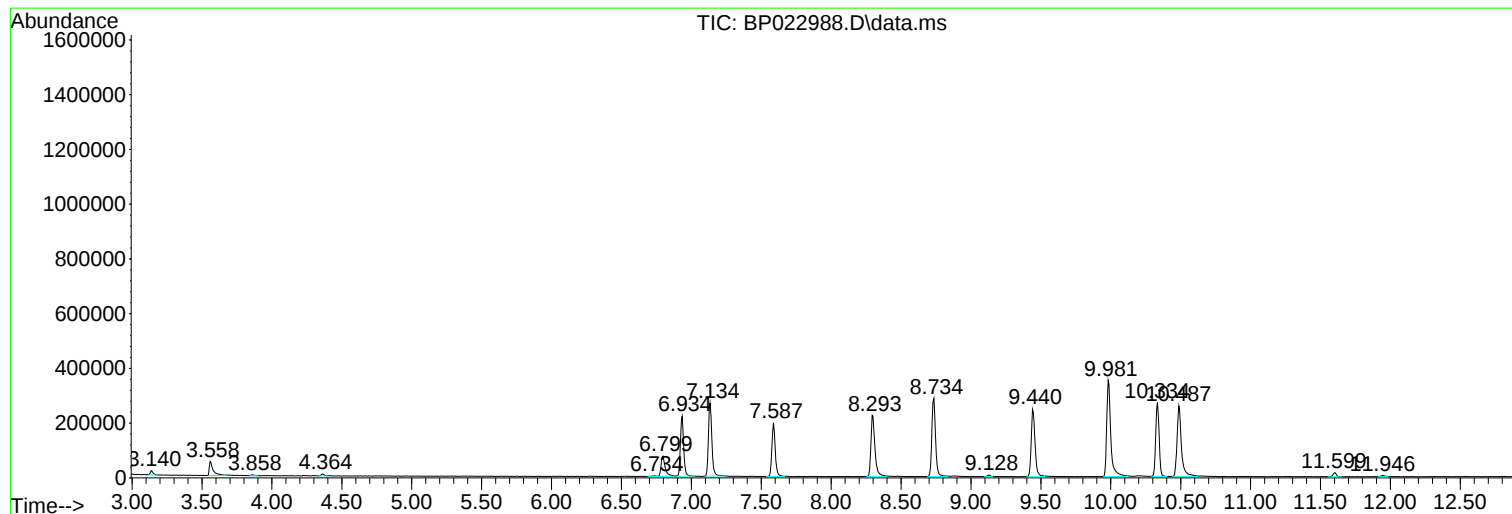
Sum of corrected areas: 22353096

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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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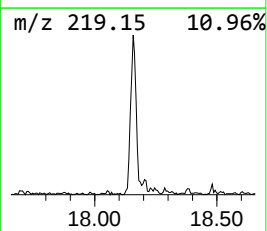
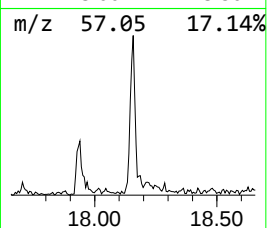
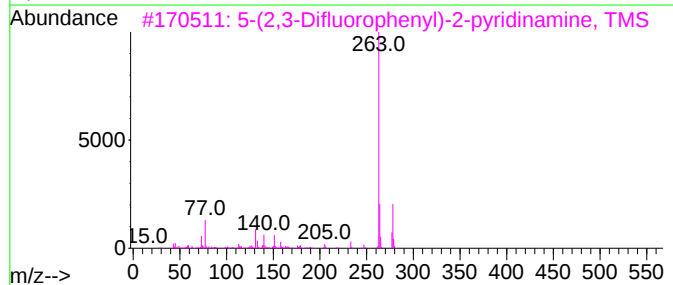
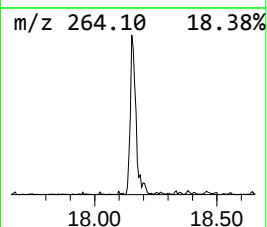
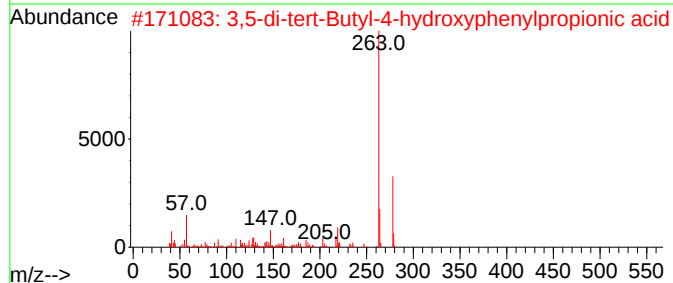
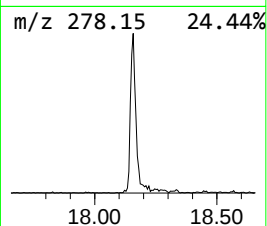
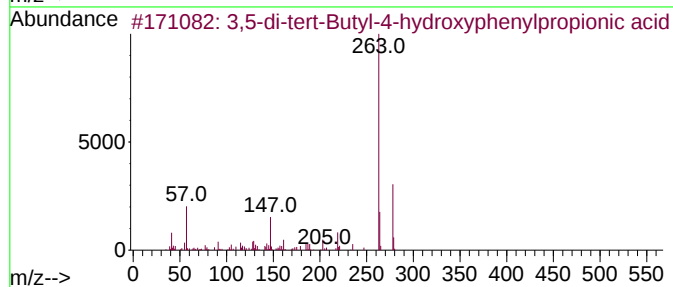
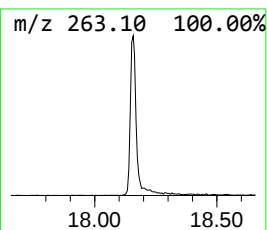
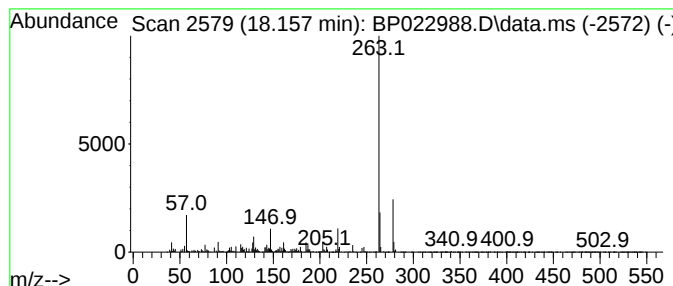
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.157	3.08 ng/ul	147577	Phenanthrene-d10	17.010

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3,5-di-tert-Butyl-4-hydroxypheny...	278	C17H26O3	020170-32-5	96
2			3,5-di-tert-Butyl-4-hydroxypheny...	278	C17H26O3	020170-32-5	94
3			5-(2,3-Difluorophenyl)-2-pyridin...	278	C14H16F2N2Si	1000504-49-4	81
4			Cyanic acid, (1-methylethylidene...	278	C17H14N2O2	001156-51-0	72
5			Cyanic acid, (1-methylethylidene...	278	C17H14N2O2	001156-51-0	64



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TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
3,5-di-tert-But...	18.157	3.1	ng/ul	147577	4	17.010	958475	20.0