

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111124\
 Data File : BG063326.D
 Acq On : 12 Nov 2024 15:17
 Operator : RC/JU
 Sample : P4801-01
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 GCPA2

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_G\Methods\SFAM-EPA-BG110624.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BG063326.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.299	32	36	43	rVB4	26379	47795	1.00%	0.114%
2	3.705	101	105	116	rBV	90965	159070	3.32%	0.380%
3	4.333	206	212	213	rBV6	4320	7458	0.16%	0.018%
4	4.539	242	247	250	rBV3	24841	35410	0.74%	0.085%
5	5.638	431	434	437	rBV3	5229	8565	0.18%	0.020%
6	5.755	450	454	457	rBV4	3629	6965	0.15%	0.017%
7	6.660	603	608	613	rVB2	6692	8385	0.17%	0.020%
8	6.819	631	635	639	rBV5	5455	9133	0.19%	0.022%
9	7.007	660	667	675	rBV	174292	304185	6.35%	0.727%
10	7.165	688	694	702	rBV	349296	549891	11.47%	1.314%
11	7.365	722	728	738	rBV	417589	694958	14.50%	1.660%
12	7.835	802	808	815	rBV	367438	617051	12.87%	1.474%
13	8.540	921	928	936	rBV2	495358	847869	17.69%	2.026%
14	8.987	997	1004	1012	rBV	506957	858413	17.91%	2.051%
15	9.051	1014	1015	1022	rVB4	5373	8469	0.18%	0.020%
16	9.269	1049	1052	1054	rBV2	7002	7082	0.15%	0.017%
17	9.421	1072	1078	1082	rBV2	8595	12628	0.26%	0.030%
18	9.709	1120	1127	1138	rVB	448639	802592	16.74%	1.918%
19	10.250	1211	1219	1231	rBV	720687	1281344	26.73%	3.061%
20	10.555	1266	1271	1273	rBV2	4246	7214	0.15%	0.017%
21	10.620	1275	1282	1289	rBV	604129	1066579	22.25%	2.548%
22	10.761	1298	1306	1314	rBV	548446	1040877	21.71%	2.487%
23	11.484	1425	1429	1431	rBV4	5192	6800	0.14%	0.016%
24	11.625	1449	1453	1455	rBV3	4286	7003	0.15%	0.017%
25	11.901	1498	1500	1507	rVB5	12769	15422	0.32%	0.037%
26	12.212	1548	1553	1559	rBV4	10065	21810	0.45%	0.052%
27	13.364	1747	1749	1753	rVB4	12437	16311	0.34%	0.039%
28	13.875	1829	1836	1844	rBV	1542234	2218773	46.28%	5.301%
29	14.163	1876	1885	1896	rVB	1598095	2494178	52.03%	5.959%
30	14.468	1930	1937	1944	rBV	1165855	1805022	37.65%	4.313%
31	14.686	1967	1974	1990	rBV3	135403	290236	6.05%	0.693%
32	15.273	2069	2074	2077	rBV5	5919	10752	0.22%	0.026%
33	15.467	2099	2107	2116	rBV	2324416	3420951	71.36%	8.174%
34	15.591	2122	2128	2137	rBV	710583	1070591	22.33%	2.558%
35	15.708	2144	2148	2153	rVB7	13891	22248	0.46%	0.053%
36	16.372	2255	2261	2264	rVB7	4747	7217	0.15%	0.017%

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37	16.695	2311	2316	2319	rBV3	6235	12213	0.25%	0.029%
38	17.148	2389	2393	2396	rVB4	33101	39227	0.82%	0.094%
39	17.224	2399	2406	2415	rVV	1562512	2315135	48.30%	5.531%
40	17.318	2416	2422	2432	rVV2	2653859	3997532	83.39%	9.551%
41	17.500	2451	2453	2456	rBV4	5685	9280	0.19%	0.022%
42	17.894	2513	2520	2523	rBV8	31757	52153	1.09%	0.125%
43	18.123	2556	2559	2566	rBV6	22373	37773	0.79%	0.090%
44	18.188	2566	2570	2573	rBV4	7496	11386	0.24%	0.027%
45	18.340	2591	2596	2604	rBV	283378	503304	10.50%	1.203%
46	18.611	2640	2642	2645	rBV4	10021	9539	0.20%	0.023%
47	18.728	2661	2662	2667	rVB5	13411	12965	0.27%	0.031%
48	19.069	2718	2720	2724	rVB4	16204	16630	0.35%	0.040%
49	19.186	2735	2740	2743	rBV2	36619	48995	1.02%	0.117%
50	19.257	2746	2752	2756	rBV	41751	67980	1.42%	0.162%
51	19.551	2800	2802	2806	rVB5	14874	15601	0.33%	0.037%
52	19.621	2806	2814	2821	rBV	3547790	4793730	100.00%	11.453%
53	20.109	2895	2897	2900	rBV4	8191	6943	0.14%	0.017%
54	20.597	2978	2980	2983	rVB4	13056	9200	0.19%	0.022%
55	20.855	3022	3024	3025	rBV2	13702	9895	0.21%	0.024%
56	21.478	3123	3130	3136	rBV	1728404	2641803	55.11%	6.312%
57	24.304	3601	3611	3625	rVB	1873073	4727325	98.61%	11.295%
58	24.510	3637	3646	3656	rVB	1031759	2726057	56.87%	6.513%

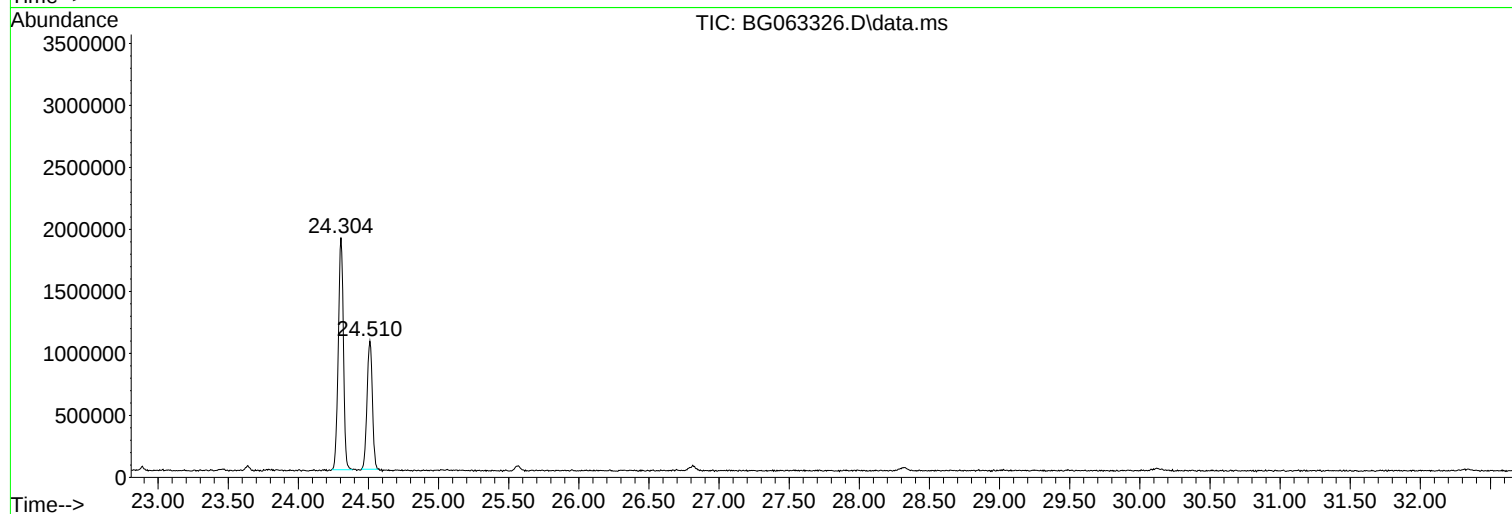
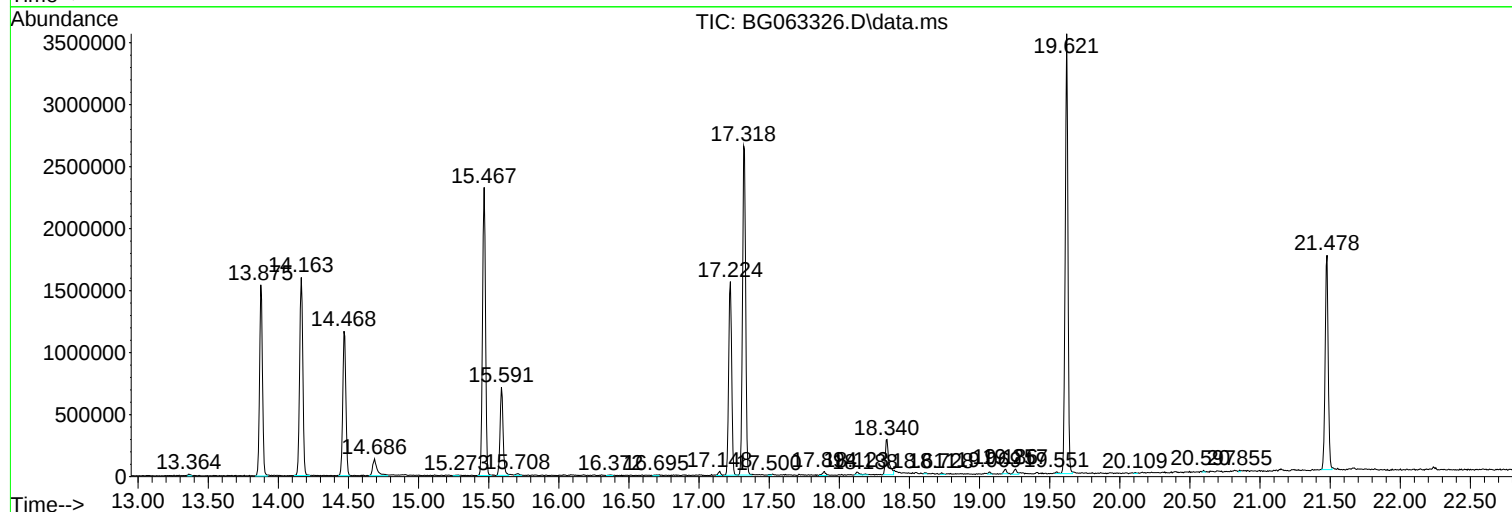
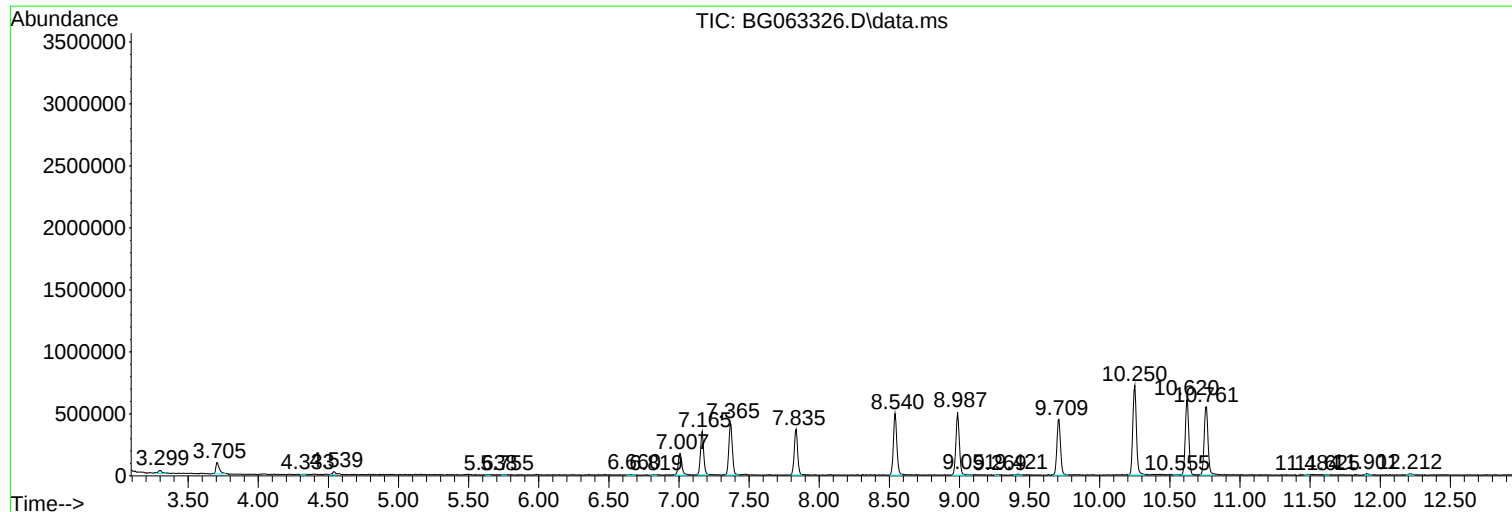
Sum of corrected areas: 41853913

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110624.MA.M
Quant Title : SVOA CALIBRATION

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



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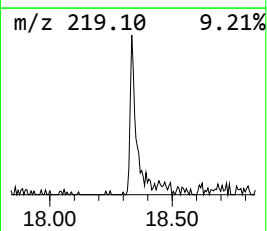
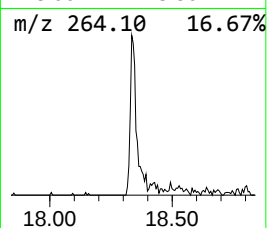
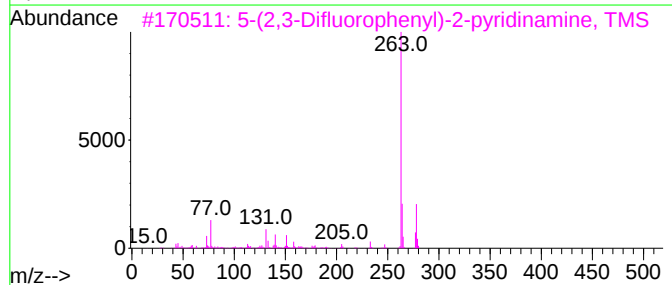
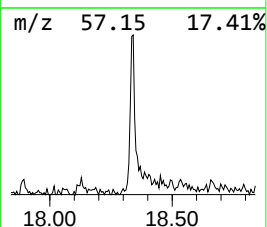
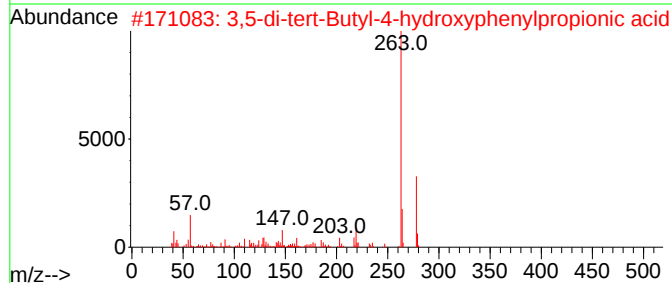
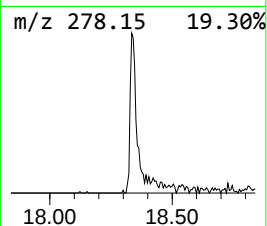
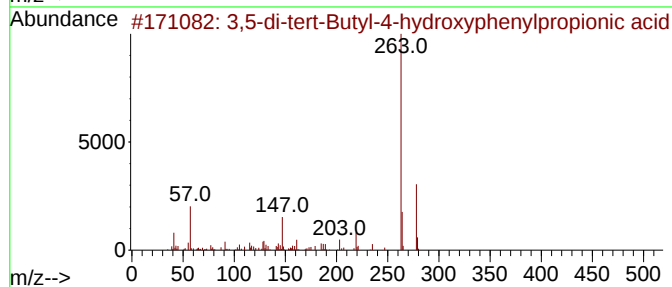
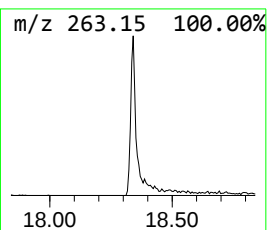
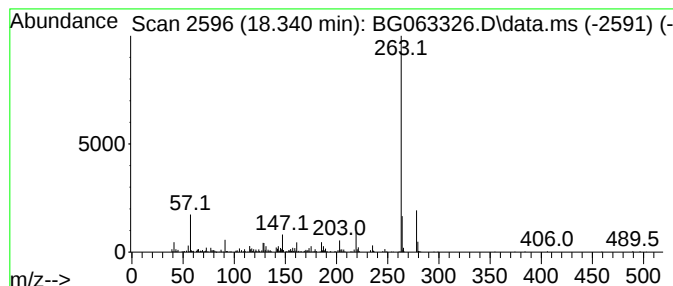
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110624.MA.M
Quant Title : SVOA CALIBRATION

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.
18.340	4.35 ng/ul	503304	Phenanthrene-d10	17.224

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3,5-di-tert-Butyl-4-hydroxypheny...	278	C17H26O3	020170-32-5	95
2			3,5-di-tert-Butyl-4-hydroxypheny...	278	C17H26O3	020170-32-5	78
3			5-(2,3-Difluorophenyl)-2-pyridin...	278	C14H16F2N2Si	1000504-49-4	72
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 Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
3,5-di-tert-But...	18.340	4.3	ng/u1	503304	4	17.224	2315140	20.0