

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG073124\
 Data File : BG062228.D
 Acq On : 31 Jul 2024 21:48
 Operator : MA/JU
 Sample : P3398-01
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 E08K3

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-BG073124.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BG062228.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.441	22	26	35	rVB2	9086	16420	0.99%	0.113%
2	3.846	91	95	104	rBV	44008	74821	4.52%	0.515%
3	4.187	147	153	164	rBV	180399	283614	17.12%	1.954%
4	4.551	213	215	218	rVB3	1839	1735	0.10%	0.012%
5	5.097	304	308	314	rBV5	1799	3072	0.19%	0.021%
6	7.124	646	653	660	rBV	49182	87604	5.29%	0.603%
7	7.306	677	684	693	rBV	137595	229479	13.85%	1.581%
8	7.506	712	718	728	rBV	148679	261129	15.76%	1.799%
9	7.617	732	737	741	rBV4	1438	2337	0.14%	0.016%
10	7.976	792	798	806	rBV	153425	262679	15.85%	1.809%
11	8.040	806	809	816	rVB5	2012	3509	0.21%	0.024%
12	8.669	909	916	924	rBV2	144660	246975	14.90%	1.701%
13	9.127	987	994	1002	rBV	157035	269388	16.26%	1.856%
14	9.303	1019	1024	1028	rBV4	1634	2367	0.14%	0.016%
15	9.850	1110	1117	1125	rBV	100103	179260	10.82%	1.235%
16	10.384	1201	1208	1218	rBV	209857	365663	22.07%	2.519%
17	10.772	1264	1274	1283	rBV	229711	416495	25.13%	2.869%
18	10.901	1287	1296	1304	rBV	210726	385095	23.24%	2.653%
19	11.553	1403	1407	1410	rBV4	1927	2754	0.17%	0.019%
20	11.712	1428	1434	1441	rBV3	1423	2893	0.17%	0.020%
21	12.352	1539	1543	1549	rVB	4516	6270	0.38%	0.043%
22	13.439	1722	1728	1734	rBV	52840	75968	4.58%	0.523%
23	14.009	1818	1825	1836	rVB	479632	685123	41.35%	4.719%
24	14.291	1866	1873	1882	rVB	504896	782540	47.22%	5.390%
25	14.596	1918	1925	1933	rVB	420855	646437	39.01%	4.453%
26	14.761	1948	1953	1960	rBV	52247	87228	5.26%	0.601%
27	15.295	2040	2044	2050	rVB2	1891	2384	0.14%	0.016%
28	15.407	2061	2063	2069	rBV3	1246	1770	0.11%	0.012%
29	15.589	2086	2094	2102	rBV	700059	1049286	63.32%	7.228%
30	15.689	2105	2111	2119	rBV	109092	157397	9.50%	1.084%
31	15.830	2131	2135	2139	rBV	2313	3085	0.19%	0.021%
32	16.223	2197	2202	2208	rVB3	1757	2973	0.18%	0.020%
33	17.339	2385	2392	2401	rVB	593698	876636	52.90%	6.039%
34	17.433	2401	2408	2417	rBV	841379	1271895	76.76%	8.761%
35	18.244	2541	2546	2550	rBV4	5144	9602	0.58%	0.066%
36	18.309	2555	2557	2562	rVB4	1713	2650	0.16%	0.018%

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Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 0.1 % of largest Peak

Max Peaks: 100

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-BG073124.MA.M
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37	19.284	2718	2723	2727	rBV4	11608	17983	1.09%	0.124%
38	19.354	2730	2735	2740	rVB	14771	19910	1.20%	0.137%
39	19.437	2748	2749	2755	rVB4	2003	2669	0.16%	0.018%
40	19.519	2759	2763	2765	rBV2	3262	2761	0.17%	0.019%
41	19.672	2786	2789	2790	rBV	4552	3754	0.23%	0.026%
42	19.719	2791	2797	2804	rVB	1134738	1593652	96.17%	10.978%
43	19.954	2833	2837	2840	rBV4	1580	2436	0.15%	0.017%
44	20.659	2954	2957	2959	rVB3	1813	1691	0.10%	0.012%
45	20.717	2963	2967	2970	rVV3	4768	4875	0.29%	0.034%
46	21.293	3062	3065	3067	rVV3	2515	2389	0.14%	0.016%
47	21.522	3101	3104	3107	rVB5	2196	2775	0.17%	0.019%
48	21.581	3107	3114	3123	rBV	625636	1038233	62.65%	7.152%
49	21.839	3152	3158	3161	rBV5	4099	6473	0.39%	0.045%
50	22.004	3184	3186	3188	rBV3	1847	2168	0.13%	0.015%
51	22.433	3256	3259	3265	rVB5	5823	9810	0.59%	0.068%
52	22.532	3275	3276	3280	rBV3	1601	1731	0.10%	0.012%
53	22.579	3282	3284	3287	rBV4	1572	1737	0.10%	0.012%
54	22.773	3313	3317	3318	rBV3	2481	3109	0.19%	0.021%
55	23.120	3372	3376	3380	rVB3	9910	14383	0.87%	0.099%
56	23.196	3386	3389	3391	rBV4	2116	2849	0.17%	0.020%
57	23.619	3459	3461	3467	rVB6	2369	3587	0.22%	0.025%
58	23.919	3505	3512	3518	rBV4	16762	34251	2.07%	0.236%
59	24.060	3534	3536	3538	rBV2	1725	2080	0.13%	0.014%
60	24.489	3596	3609	3621	rBV	606598	1657083	100.00%	11.415%
61	24.700	3633	3645	3657	rVB	400114	1102902	66.56%	7.597%
62	24.794	3660	3661	3664	rBV3	1931	2115	0.13%	0.015%
63	24.853	3664	3671	3680	rVB3	21391	54081	3.26%	0.373%
64	25.241	3733	3737	3738	rBV3	1736	2318	0.14%	0.016%
65	25.675	3808	3811	3813	rBV4	2393	2877	0.17%	0.020%
66	25.863	3841	3843	3844	rBV2	2163	1809	0.11%	0.012%
67	25.969	3851	3861	3870	rBV3	20926	61745	3.73%	0.425%
68	27.320	4080	4091	4101	rBV4	17820	65822	3.97%	0.453%
69	27.561	4129	4132	4133	rBV3	1816	1684	0.10%	0.012%
70	28.248	4247	4249	4251	rVB3	2120	1994	0.12%	0.014%
71	28.272	4251	4253	4254	rBV2	1718	1723	0.10%	0.012%
72	28.924	4355	4364	4365	rBV4	11011	21119	1.27%	0.145%
73	29.529	4465	4467	4470	rBV4	1715	1890	0.11%	0.013%
74	30.216	4583	4584	4586	rBV2	2682	2066	0.12%	0.014%
75	30.240	4586	4588	4590	rVB2	2303	2034	0.12%	0.014%

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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_G\Methods\SFAM-EPA-BG073124.MA.M
Title : SVOA CALIBRATION

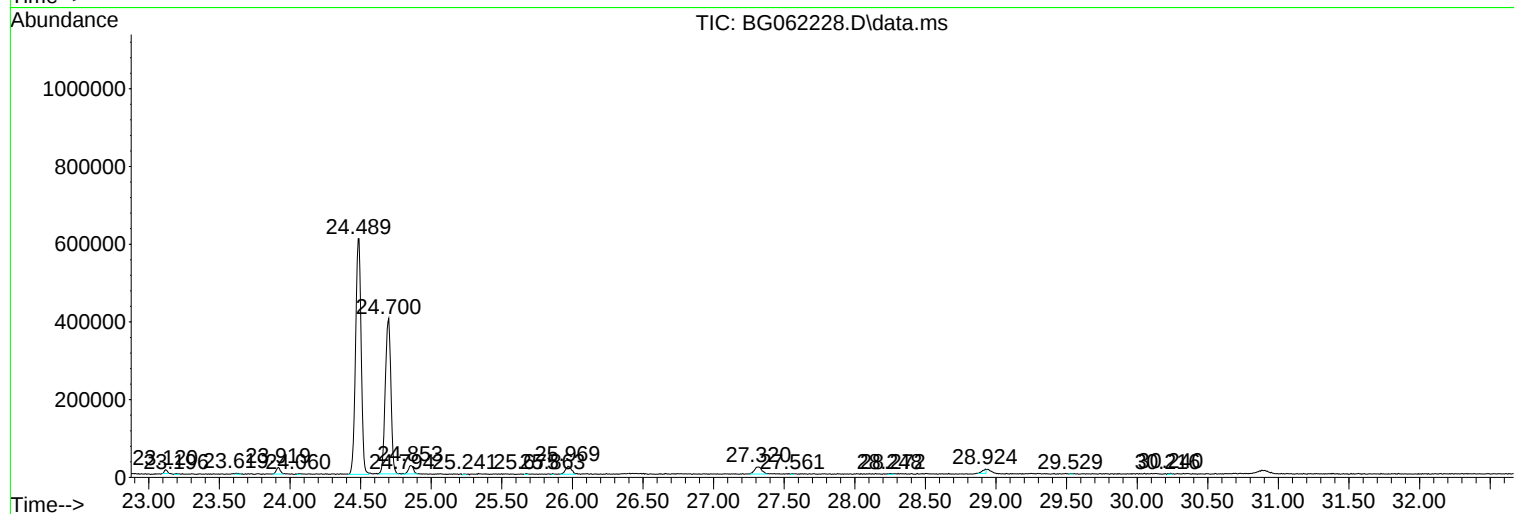
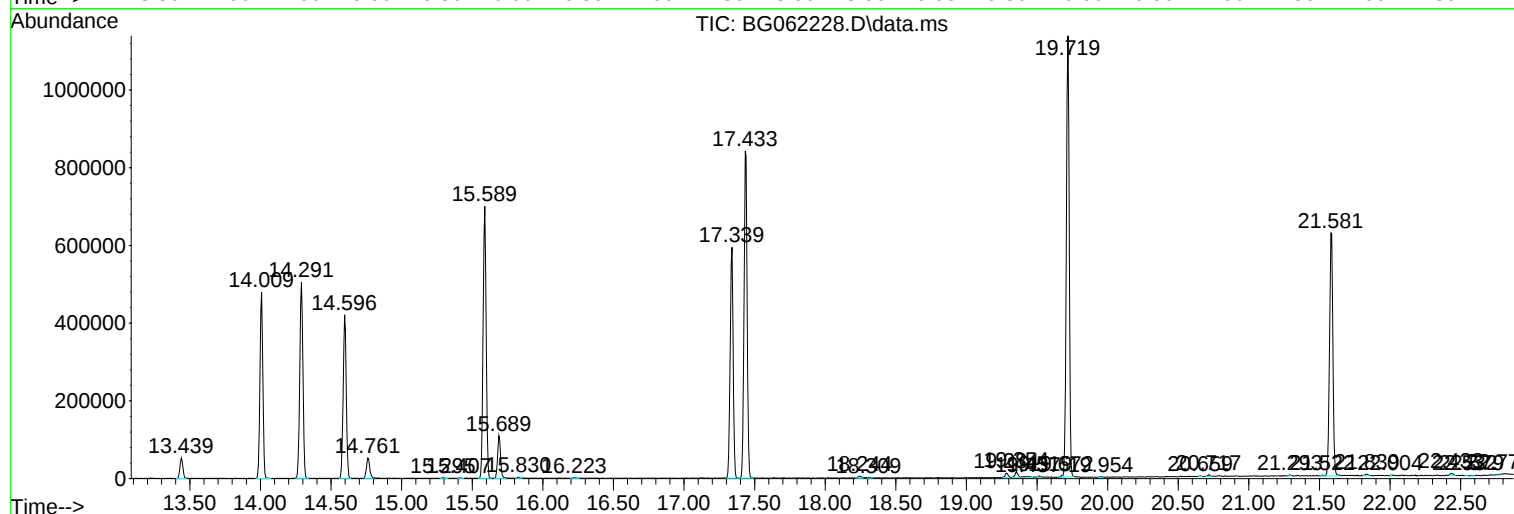
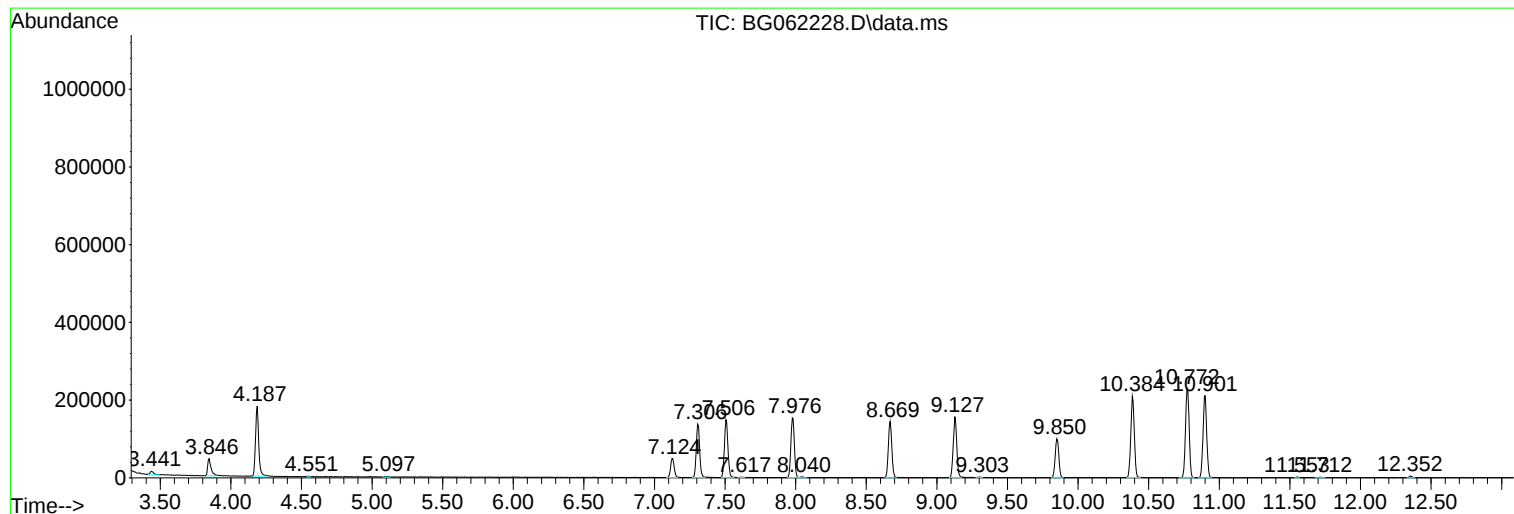
Sum of corrected areas: 14517101

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG073124.MA.M
Quant Title : SVOA CALIBRATION

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



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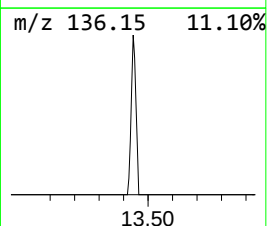
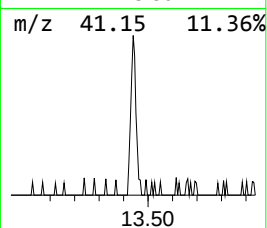
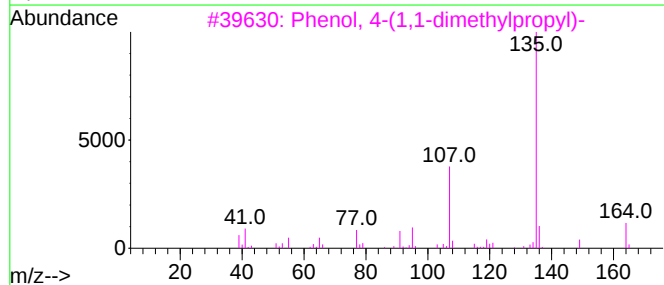
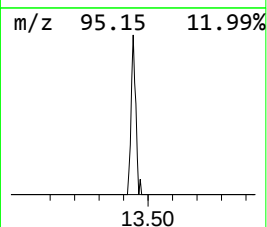
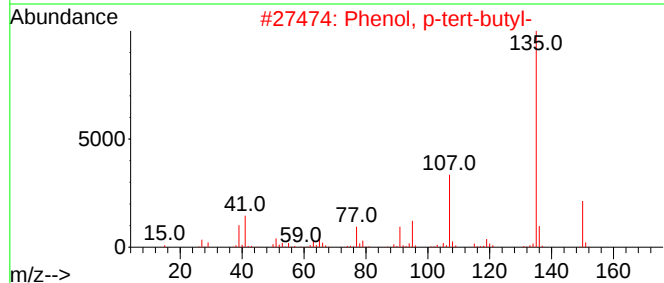
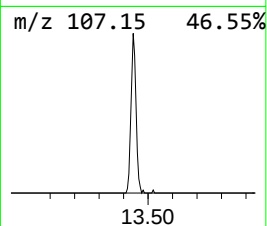
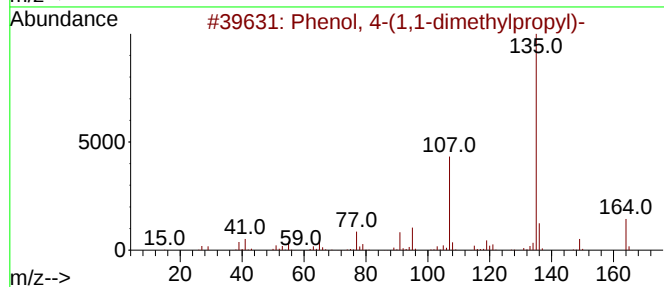
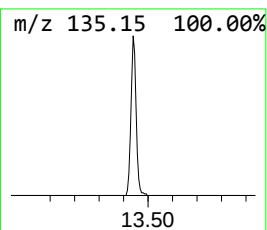
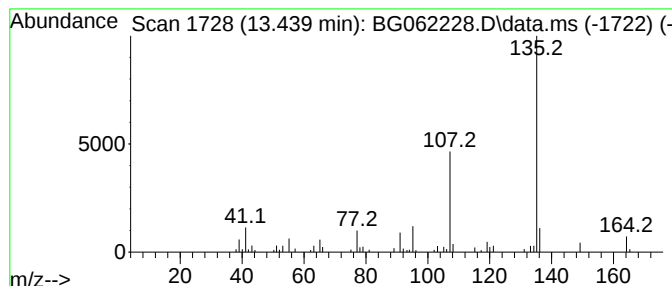
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG073124.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Phenol, 4-(1,1-dimethylprop... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.439	2.35 ng/ul	75968	Acenaphthene-d10	14.596

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	93
2			Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	90
3			Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	87
4			Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	72
5			Hexestrol, O-trifluoroacetyl-	366	C20H21F3O3	1000365-31-7	72



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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Phenol, 4-(1,1-...	13.439	2.4	ng/ul	75968	3	14.596	646437	20.0