

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG010923\
 Data File : BG056230.D
 Acq On : 9 Jan 2023 11:43
 Operator : CG/JU
 Sample : PB150060BL
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK060

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-BG010423.M
 Title : SVOA CALIBRATION

Signal : TIC: BG056230.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.353	6	11	18	rVB	19542	28465	1.75%	0.186%
2	3.541	38	43	46	rBV3	1231	2210	0.14%	0.014%
3	3.629	53	58	64	rBV3	9221	14979	0.92%	0.098%
4	4.058	128	131	144	rBV	109006	178848	10.99%	1.169%
5	4.416	185	192	197	rBV	8367	13188	0.81%	0.086%
6	7.460	700	710	720	rBB	158415	273700	16.82%	1.789%
7	7.636	731	740	748	rBB	133303	221099	13.59%	1.445%
8	7.854	769	777	787	rBV	158034	275432	16.93%	1.801%
9	8.330	852	858	865	rVB	158853	255159	15.68%	1.668%
10	9.017	968	975	984	rBV	228519	385542	23.70%	2.520%
11	9.499	1049	1057	1066	rBB	151343	269241	16.55%	1.760%
12	10.227	1175	1181	1189	rBV	130186	235321	14.46%	1.538%
13	10.768	1266	1273	1283	rBV	240959	428411	26.33%	2.801%
14	11.162	1332	1340	1347	rBV	211832	384512	23.63%	2.514%
15	11.291	1353	1362	1371	rBV	208643	385614	23.70%	2.521%
16	12.718	1600	1605	1612	rBV4	4110	7834	0.48%	0.051%
17	14.328	1873	1879	1886	rBV	516891	739771	45.47%	4.836%
18	14.640	1925	1932	1940	rBV	577145	904825	55.61%	5.915%
19	14.945	1973	1984	1991	rBV	430074	686038	42.17%	4.485%
20	15.104	2005	2011	2026	rBV	162057	267418	16.44%	1.748%
21	15.927	2144	2151	2159	rBV	807348	1222900	75.17%	7.994%
22	16.032	2162	2169	2175	rVV	187486	271536	16.69%	1.775%
23	16.073	2175	2176	2182	rVB	1589	1790	0.11%	0.012%
24	16.167	2188	2192	2195	rBV4	2827	4100	0.25%	0.027%
25	17.354	2389	2394	2396	rBV3	1900	2408	0.15%	0.016%
26	17.472	2408	2414	2417	rBV2	1462	2496	0.15%	0.016%
27	17.677	2439	2449	2459	rVB	582087	907934	55.81%	5.935%
28	17.777	2459	2466	2473	rBV	879624	1349277	82.93%	8.820%
29	18.048	2508	2512	2515	rVB4	1874	1875	0.12%	0.012%
30	18.224	2539	2542	2544	rBV2	1871	1867	0.11%	0.012%
31	18.318	2556	2558	2560	rVB2	2133	1871	0.12%	0.012%
32	18.664	2613	2617	2619	rBV4	1535	2157	0.13%	0.014%
33	19.246	2714	2716	2720	rBV4	2005	2508	0.15%	0.016%
34	19.528	2761	2764	2765	rBV3	1515	1696	0.10%	0.011%
35	19.610	2774	2778	2784	rVB4	14372	21852	1.34%	0.143%
36	19.687	2784	2791	2796	rBV2	11943	22875	1.41%	0.150%

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Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG010423.M
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37	19.798	2807	2810	2812	rBV2	1380	1628	0.10%	0.011%
38	19.922	2829	2831	2833	rBV	1789	1819	0.11%	0.012%
39	19.992	2841	2843	2845	rBV3	2301	2146	0.13%	0.014%
40	20.045	2845	2852	2859	rBV	1147965	1626947	100.00%	10.636%
41	20.157	2869	2871	2874	rBV3	1938	2182	0.13%	0.014%
42	20.433	2917	2918	2920	rBV	1879	1684	0.10%	0.011%
43	20.627	2949	2951	2953	rVB3	4255	2667	0.16%	0.017%
44	20.739	2966	2970	2971	rBV4	3683	4548	0.28%	0.030%
45	20.968	3006	3009	3011	rVB4	4232	4844	0.30%	0.032%
46	21.026	3017	3019	3021	rBV3	3251	3373	0.21%	0.022%
47	21.209	3048	3050	3051	rBV2	3136	2490	0.15%	0.016%
48	21.496	3096	3099	3100	rBV3	4540	3758	0.23%	0.025%
49	21.978	3174	3181	3191	rVB2	593300	1006250	61.85%	6.578%
50	22.783	3316	3318	3320	rBV3	4893	2220	0.14%	0.015%
51	23.523	3441	3444	3450	rVB8	10180	18585	1.14%	0.121%
52	24.405	3588	3594	3599	rVB7	14339	27522	1.69%	0.180%
53	24.916	3680	3681	3683	rBV2	4512	3170	0.19%	0.021%
54	25.204	3717	3730	3744	rVB2	551377	1607977	98.83%	10.512%
55	25.439	3758	3770	3781	rVB3	354452	1054518	64.82%	6.893%
56	25.756	3822	3824	3825	rBV2	3960	2787	0.17%	0.018%
57	26.068	3875	3877	3879	rVB3	3251	2004	0.12%	0.013%
58	26.449	3940	3942	3944	rVB3	3871	2261	0.14%	0.015%
59	26.655	3969	3977	3986	rVB9	19956	64468	3.96%	0.421%
60	26.867	4011	4013	4015	rBV3	3401	2385	0.15%	0.016%
61	27.096	4049	4052	4053	rBV3	3633	3375	0.21%	0.022%
62	28.012	4206	4208	4210	rBV3	4820	4756	0.29%	0.031%
63	28.259	4248	4250	4251	rBV2	5276	4461	0.27%	0.029%
64	29.082	4388	4390	4393	rBV4	3595	3202	0.20%	0.021%
65	29.140	4398	4400	4402	rBV3	4934	4937	0.30%	0.032%
66	29.851	4518	4521	4523	rBV3	7980	10961	0.67%	0.072%
67	29.898	4524	4529	4530	rVV4	9674	13834	0.85%	0.090%
68	30.022	4548	4550	4552	rBV3	4140	4186	0.26%	0.027%
69	30.151	4570	4572	4573	rVB2	5634	2706	0.17%	0.018%
70	30.404	4613	4615	4619	rVB5	4846	5355	0.33%	0.035%
71	30.433	4619	4620	4621	rBV	5242	2444	0.15%	0.016%
72	31.244	4756	4758	4760	rBV3	4070	2473	0.15%	0.016%
73	32.542	4976	4979	4982	rBV5	4452	5638	0.35%	0.037%

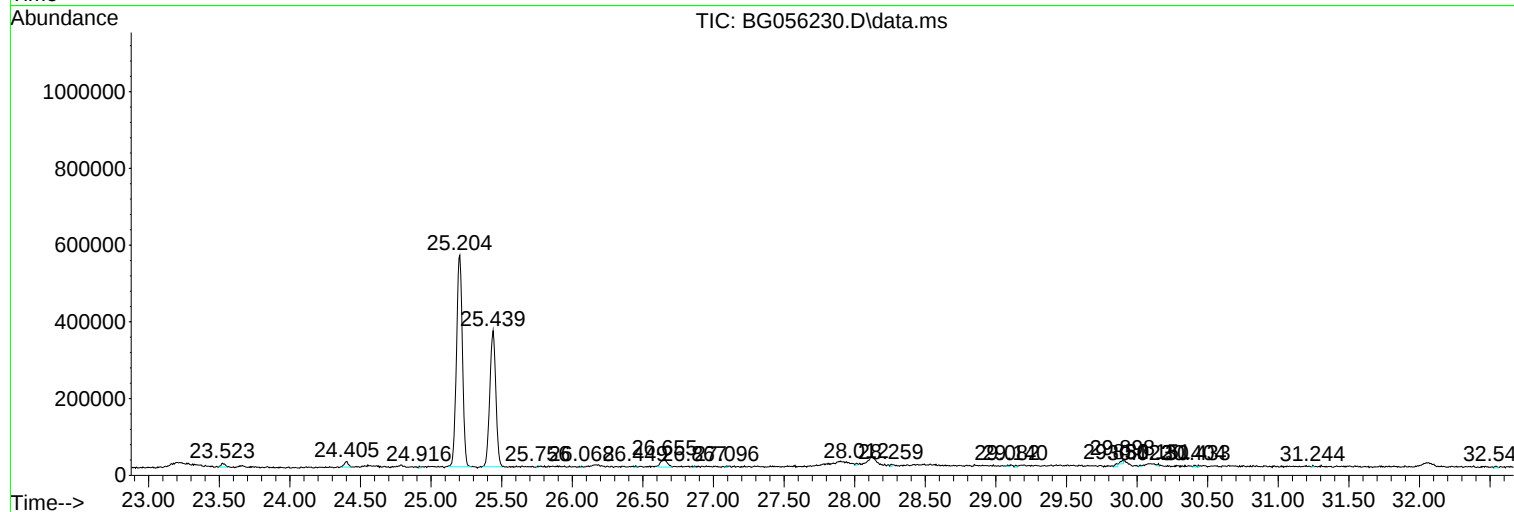
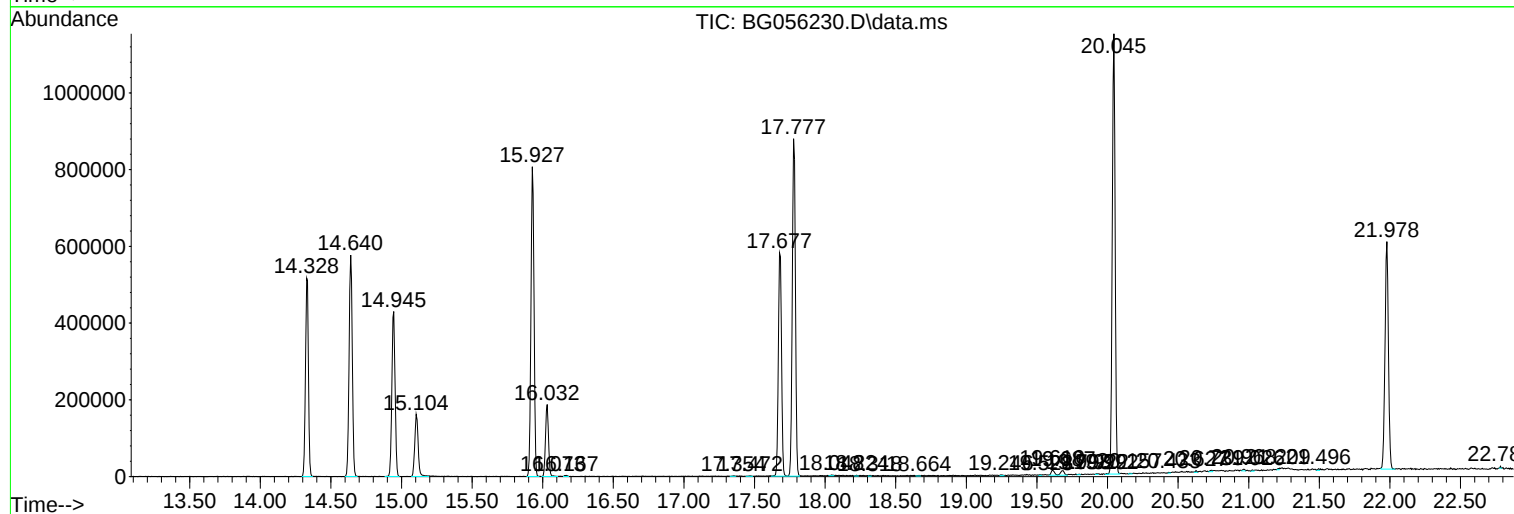
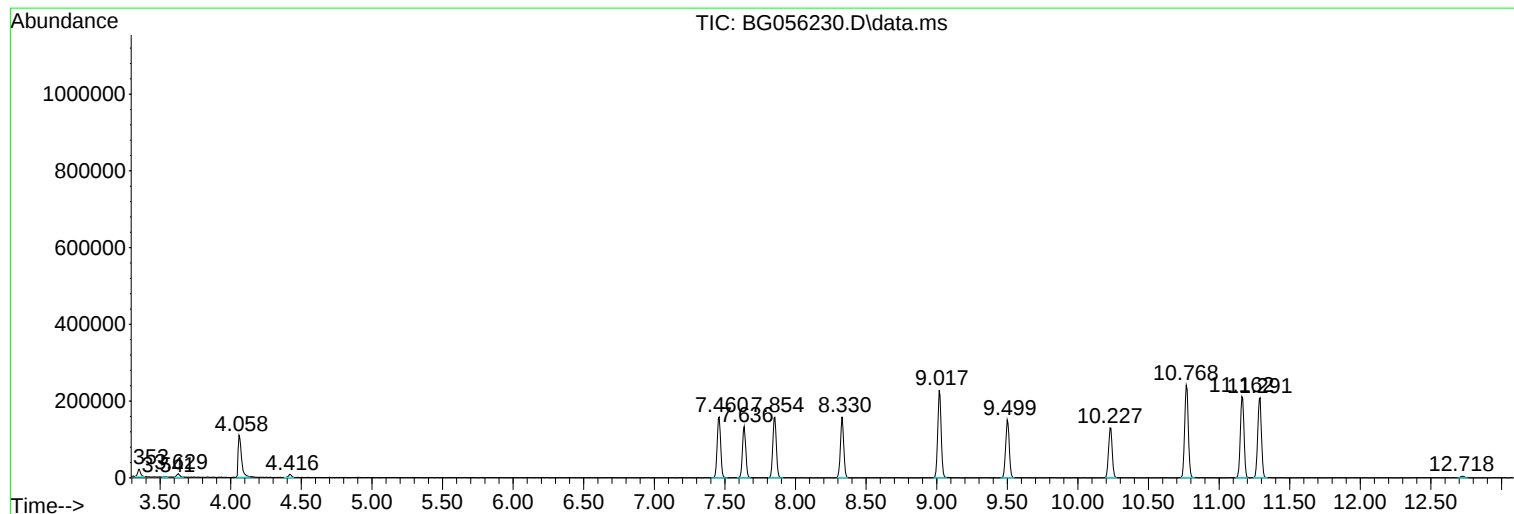
Sum of corrected areas: 15297310

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG010423.M
Quant Title : SVOA CALIBRATION

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



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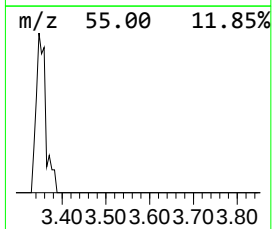
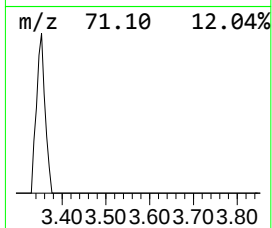
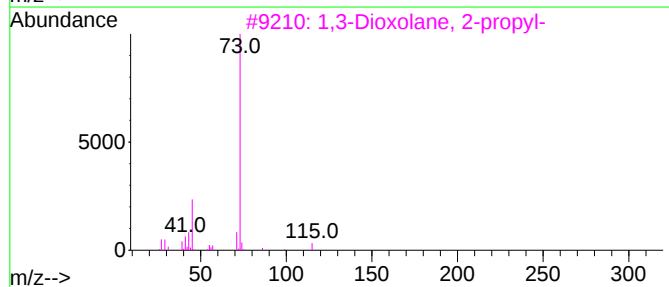
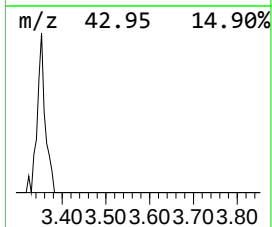
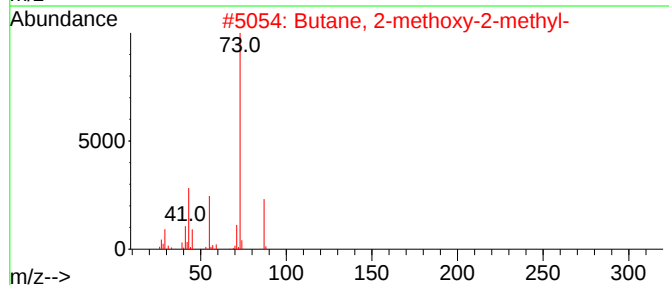
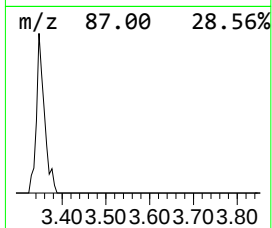
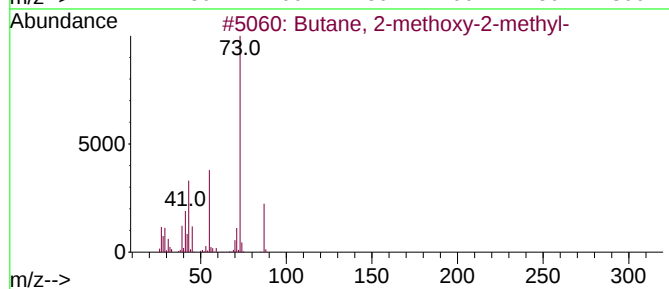
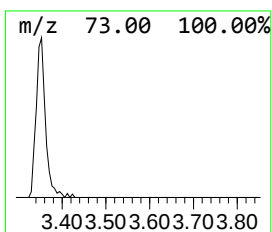
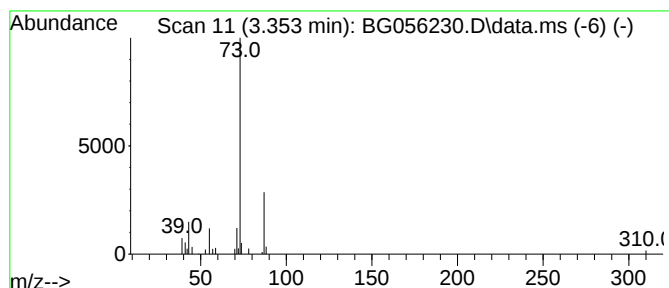
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG010423.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 unknown-01 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.353	2.23 ng/ul	28465	1,4-Dichlorobenzene-d4	8.330

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	39		
2	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	39		
3	1,3-Dioxolane, 2-propyl-	116	C6H12O2	003390-13-4	23		
4	1-Pentanamine	87	C5H13N	000110-58-7	9		
5	N-Ethylformamide	73	C3H7NO	000627-45-2	9		



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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
unknown-01	3.353	2.2	ng/u1	28465	1	8.330	255159	20.0