

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG081921\
 Data File : BG049896.D
 Acq On : 19 Aug 2021 15:15
 Operator : CG/JU
 Sample : PB138362BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK362

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-BG081921.M
 Title : SVOA CALIBRATION

Signal : TIC: BG049896.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.073	3	6	13	rVB2	15030	22265	2.98%	0.293%
2	3.232	31	33	38	rVB4	1836	2200	0.29%	0.029%
3	3.290	41	43	46	rVB3	1510	1221	0.16%	0.016%
4	3.343	49	52	58	rVB3	6492	9766	1.31%	0.128%
5	3.431	65	67	70	rBV3	1307	1063	0.14%	0.014%
6	3.531	81	84	86	rBV4	1342	1726	0.23%	0.023%
7	3.578	90	92	95	rVB	2059	1477	0.20%	0.019%
8	3.678	104	109	110	rVB2	1286	1760	0.24%	0.023%
9	3.772	121	125	139	rBV	62200	127434	17.07%	1.675%
10	4.119	181	184	185	rVB2	1624	1231	0.16%	0.016%
11	4.730	286	288	294	rVB3	744	1106	0.15%	0.015%
12	5.241	374	375	380	rBV2	1853	1901	0.25%	0.025%
13	5.364	393	396	399	rBV2	916	1430	0.19%	0.019%
14	5.658	444	446	448	rBV2	1258	1228	0.16%	0.016%
15	5.799	468	470	473	rVB	1268	1115	0.15%	0.015%
16	5.934	491	493	495	rVB	1629	1110	0.15%	0.015%
17	6.116	522	524	530	rBV3	1080	1618	0.22%	0.021%
18	6.298	552	555	558	rBV	1230	1473	0.20%	0.019%
19	6.363	564	566	568	rVB2	1120	1041	0.14%	0.014%
20	6.856	647	650	653	rBV2	1591	1406	0.19%	0.018%
21	7.144	692	699	711	rBV	104265	194737	26.08%	2.560%
22	7.303	720	726	734	rVB	72965	124044	16.61%	1.631%
23	7.391	739	741	744	rBV2	1120	1509	0.20%	0.020%
24	7.514	756	762	772	rVB	110628	195418	26.17%	2.569%
25	7.837	815	817	821	rVB	1135	1224	0.16%	0.016%
26	7.878	821	824	828	rBV2	948	1616	0.22%	0.021%
27	7.984	836	842	848	rVB2	85478	142607	19.10%	1.875%
28	8.078	857	858	862	rBV2	1180	1139	0.15%	0.015%
29	8.107	862	863	867	rVB	1087	1111	0.15%	0.015%
30	8.254	885	888	890	rBV	1155	1238	0.17%	0.016%
31	8.701	957	964	972	rBV	126563	233639	31.29%	3.071%
32	9.153	1033	1041	1049	rBV2	111762	198934	26.64%	2.615%
33	9.417	1085	1086	1089	rBV	1098	1172	0.16%	0.015%
34	9.605	1116	1118	1123	rVB	1346	1469	0.20%	0.019%
35	9.652	1123	1126	1127	rBV	1462	1186	0.16%	0.016%
36	9.881	1159	1165	1174	rVB	98773	182999	24.51%	2.406%

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Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG081921.M
 Title : SVOA CALIBRATION

37	10.011	1184	1187	1190	rBV2	858	1105	0.15%	0.015%
38	10.163	1209	1213	1216	rBV	936	1161	0.16%	0.015%
39	10.428	1252	1258	1267	rBV	131225	244461	32.74%	3.214%
40	10.804	1315	1322	1329	rBV	117285	206381	27.64%	2.713%
41	10.939	1338	1345	1357	rBV	128014	240896	32.26%	3.167%
42	11.191	1387	1388	1393	rBV	684	1052	0.14%	0.014%
43	11.315	1407	1409	1413	rVB	1060	1151	0.15%	0.015%
44	12.390	1588	1592	1594	rBV2	1749	2347	0.31%	0.031%
45	14.041	1867	1873	1879	rBV	285459	426072	57.06%	5.601%
46	14.334	1915	1923	1931	rBV	301342	453639	60.75%	5.964%
47	14.640	1966	1975	1982	rBV2	198548	308713	41.34%	4.058%
48	14.857	2006	2012	2025	rBV	87423	176179	23.59%	2.316%
49	15.633	2138	2144	2155	rVB	393032	574618	76.96%	7.554%
50	15.762	2160	2166	2175	rBV2	125038	198743	26.62%	2.613%
51	15.873	2180	2185	2195	rVB3	4793	9032	1.21%	0.119%
52	16.061	2214	2217	2221	rBV2	969	1341	0.18%	0.018%
53	16.414	2275	2277	2281	rBV	1126	1303	0.17%	0.017%
54	17.283	2423	2425	2428	rBV	705	1089	0.15%	0.014%
55	17.389	2437	2443	2449	rBV	253158	385441	51.62%	5.067%
56	17.489	2453	2460	2468	rBV	425860	625631	83.79%	8.225%
57	17.706	2494	2497	2501	rVB	1291	1627	0.22%	0.021%
58	17.994	2542	2546	2549	rBV3	953	1315	0.18%	0.017%
59	19.057	2724	2727	2729	rVB3	1272	1056	0.14%	0.014%
60	19.157	2742	2744	2745	rBV	1888	1195	0.16%	0.016%
61	19.204	2750	2752	2755	rBV2	1507	1695	0.23%	0.022%
62	19.310	2768	2770	2774	rVB2	1415	1380	0.18%	0.018%
63	19.363	2777	2779	2780	rVB2	2682	1361	0.18%	0.018%
64	19.416	2784	2788	2793	rVB	6472	8367	1.12%	0.110%
65	19.457	2793	2795	2798	rBV2	1309	1166	0.16%	0.015%
66	19.651	2827	2828	2829	rBV	2302	1476	0.20%	0.019%
67	19.668	2829	2831	2837	rVB2	4289	7325	0.98%	0.096%
68	19.780	2844	2850	2857	rBV	530811	746682	100.00%	9.816%
69	19.991	2884	2886	2889	rVB2	1563	1652	0.22%	0.022%
70	20.032	2889	2893	2896	rBV5	2655	4727	0.63%	0.062%
71	20.103	2902	2905	2907	rBV4	3409	4010	0.54%	0.053%
72	20.250	2929	2930	2932	rBV2	3290	1776	0.24%	0.023%
73	20.267	2932	2933	2935	rVV2	1850	1100	0.15%	0.014%
74	20.626	2993	2994	2996	rBV2	3748	2297	0.31%	0.030%
75	20.726	3009	3011	3014	rVB4	5173	4159	0.56%	0.055%
76	21.660	3165	3170	3179	rVB	254814	407921	54.63%	5.363%

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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-BG081921.M
Title : SVOA CALIBRATION

77	24.673	3674	3683	3693	rVB	249312	657648	88.08%	8.646%
78	24.896	3711	3721	3730	rVB2	149836	412006	55.18%	5.416%
79	28.298	4298	4300	4302	rVB2	5083	2457	0.33%	0.032%
80	29.202	4452	4454	4458	rVB5	6957	4315	0.58%	0.057%

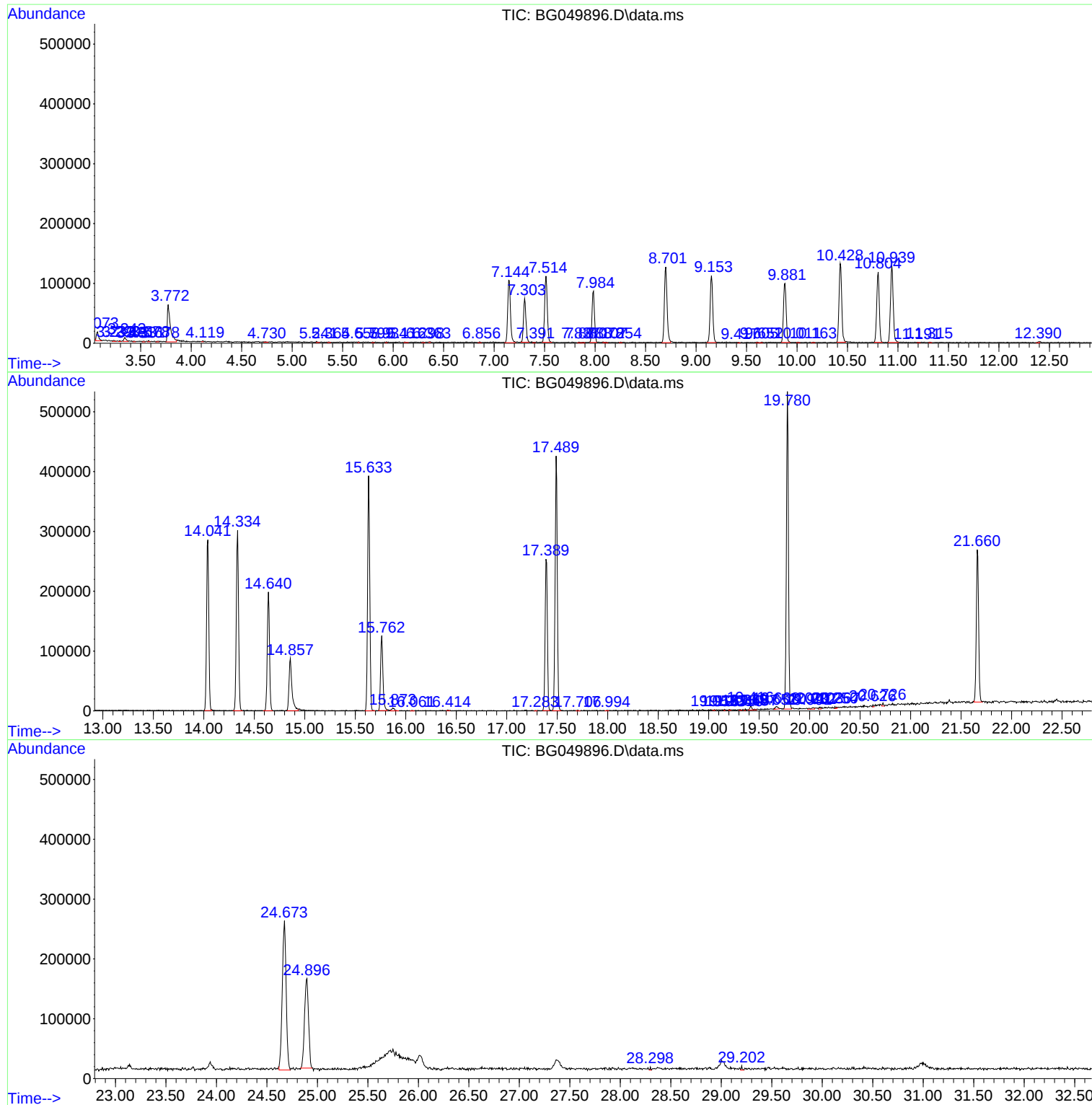
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Instrument :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG081921.M
Quant Title : SVOA CALIBRATION

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



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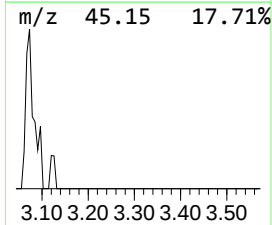
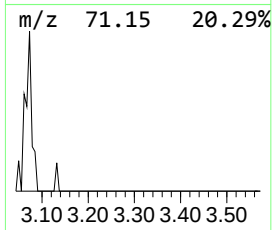
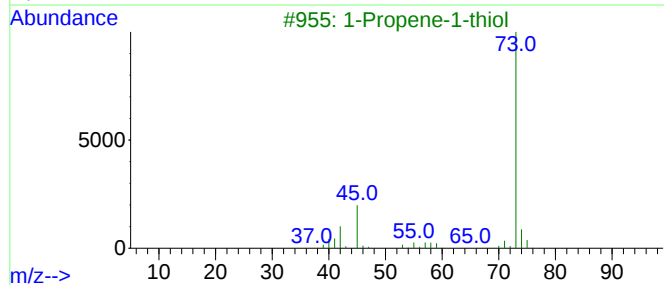
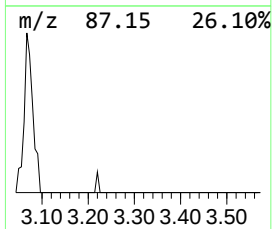
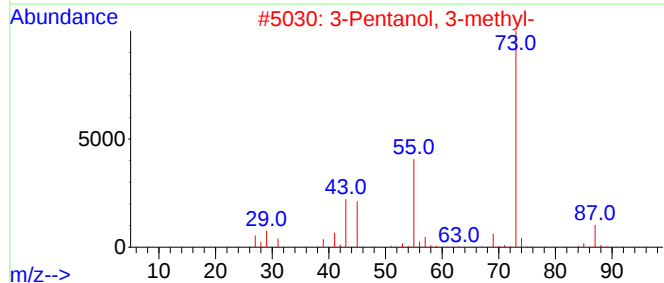
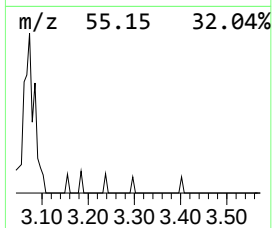
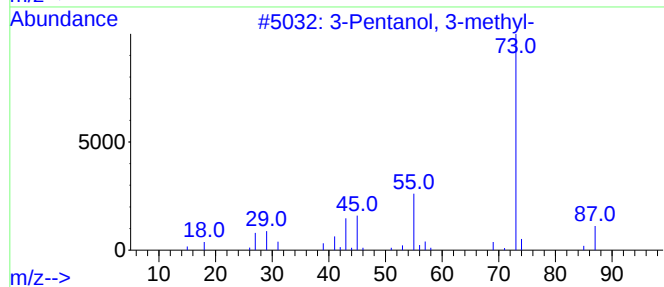
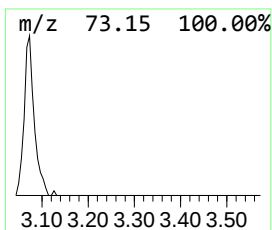
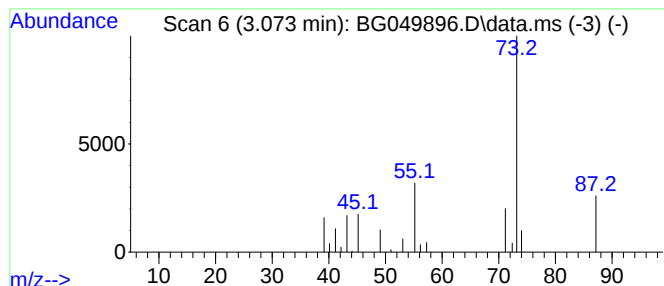
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG081921.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.073	3.12 ng/ul	22265	1,4-Dichlorobenzene-d4	7.984

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		3-Pentanol, 3-methyl-	102	C6H14O	000077-74-7	38
2		3-Pentanol, 3-methyl-	102	C6H14O	000077-74-7	17
3		1-Propene-1-thiol	74	C3H6S	000925-89-3	9
4		Ethyl formate	74	C3H6O2	000109-94-4	7
5		Methane, isothiocyanato-	73	C2H3NS	000556-61-6	7



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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
(DEL) Alkane: S...	3.073	3.1	ng/ul	22265	1	7.984	142607	20.0