

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG011422\
 Data File : BG051974.D
 Acq On : 14 Jan 2022 13:35
 Operator : CG/JU
 Sample : N1100-18
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BGLT2

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-BG010622.M
 Title : SVOA CALIBRATION

Signal : TIC: BG051974.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.570	10	14	19	rVB4	5420	8846	0.98%	0.108%
2	3.852	60	62	64	rBV2	1484	1351	0.15%	0.016%
3	3.999	83	87	96	rBV	62689	105827	11.68%	1.286%
4	4.246	126	129	133	rVB	1310	1466	0.16%	0.018%
5	4.352	144	147	151	rVB3	2070	2639	0.29%	0.032%
6	5.068	263	269	273	rVB2	900	1399	0.15%	0.017%
7	5.286	305	306	310	rVB2	2460	2146	0.24%	0.026%
8	6.578	521	526	528	rBV	812	1311	0.14%	0.016%
9	7.060	604	608	609	rVB2	1156	1177	0.13%	0.014%
10	7.383	656	663	674	rBB	99523	180621	19.94%	2.195%
11	7.547	682	691	700	rVB	64332	110070	12.15%	1.338%
12	7.770	722	729	736	rBV	85226	145167	16.03%	1.764%
13	7.923	752	755	758	rBV	1009	1176	0.13%	0.014%
14	8.246	803	810	815	rBV	76122	125324	13.83%	1.523%
15	8.293	815	818	827	rVB2	9802	14413	1.59%	0.175%
16	8.470	843	848	851	rVB	825	1447	0.16%	0.018%
17	8.945	921	929	938	rBV	119621	199805	22.06%	2.428%
18	9.415	1002	1009	1017	rVB	93946	166152	18.34%	2.019%
19	9.838	1076	1081	1089	rVB4	4453	6657	0.73%	0.081%
20	10.144	1126	1133	1141	rBV2	78692	144110	15.91%	1.752%
21	10.696	1220	1227	1239	rVB	115396	208012	22.96%	2.528%
22	10.837	1246	1251	1257	rVB3	3218	5885	0.65%	0.072%
23	11.084	1286	1293	1301	rBV	107164	186692	20.61%	2.269%
24	11.207	1307	1314	1325	rBV	123739	223694	24.69%	2.719%
25	12.476	1525	1530	1534	rBV4	2200	3427	0.38%	0.042%
26	12.658	1557	1561	1562	rBV2	1621	1928	0.21%	0.023%
27	13.281	1661	1667	1670	rBV2	1421	2548	0.28%	0.031%
28	13.422	1687	1691	1694	rVB2	1471	1795	0.20%	0.022%
29	13.698	1735	1738	1743	rBV2	4449	6569	0.73%	0.080%
30	13.833	1758	1761	1765	rVB2	973	1256	0.14%	0.015%
31	14.091	1800	1805	1812	rBV	40161	58209	6.43%	0.707%
32	14.279	1830	1837	1843	rBV	270027	390172	43.07%	4.742%
33	14.397	1854	1857	1860	rVB2	1092	1325	0.15%	0.016%
34	14.514	1873	1877	1878	rBV	1786	2138	0.24%	0.026%
35	14.585	1882	1889	1896	rVB	285711	444123	49.03%	5.398%
36	14.761	1916	1919	1924	rVB2	2839	3849	0.42%	0.047%

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Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG010622.M
 Title : SVOA CALIBRATION

37	14.890	1934	1941	1948	rVB2	192587	306183	33.80%	3.721%
38	15.061	1961	1970	1983	rBV	130852	226117	24.96%	2.748%
39	15.290	2005	2009	2011	rVB2	962	1295	0.14%	0.016%
40	15.513	2044	2047	2052	rVB	1057	1574	0.17%	0.019%

41	15.713	2078	2081	2085	rVB4	3040	3785	0.42%	0.046%
42	15.877	2102	2109	2116	rVB	386853	591755	65.33%	7.192%
43	15.983	2120	2127	2136	rBV	116907	176652	19.50%	2.147%
44	16.118	2146	2150	2154	rVB3	1888	2524	0.28%	0.031%
45	16.271	2172	2176	2179	rVB2	1814	2035	0.22%	0.025%

46	16.506	2213	2216	2220	rBV3	2153	2682	0.30%	0.033%
47	16.559	2220	2225	2226	rVV2	2327	3503	0.39%	0.043%
48	16.600	2230	2232	2238	rVB4	1596	2516	0.28%	0.031%
49	16.752	2254	2258	2263	rBV2	1881	3649	0.40%	0.044%
50	16.917	2282	2286	2287	rBV	1671	1280	0.14%	0.016%

51	17.023	2301	2304	2309	rVB	1614	1637	0.18%	0.020%
52	17.070	2309	2312	2314	rBV	1008	1387	0.15%	0.017%
53	17.217	2335	2337	2340	rBV3	1427	1900	0.21%	0.023%
54	17.310	2348	2353	2354	rBV3	1526	2248	0.25%	0.027%
55	17.369	2360	2363	2367	rVB3	4412	5071	0.56%	0.062%

56	17.416	2367	2371	2377	rVB4	2698	5597	0.62%	0.068%
57	17.469	2377	2380	2382	rBV2	1114	1364	0.15%	0.017%
58	17.528	2387	2390	2394	rVB4	2294	3289	0.36%	0.040%
59	17.634	2402	2408	2415	rBV	298548	461206	50.91%	5.606%
60	17.739	2418	2426	2433	rVV	475315	736569	81.31%	8.952%

61	17.810	2434	2438	2442	rVB6	6932	8437	0.93%	0.103%
62	18.057	2478	2480	2484	rBV2	1168	1213	0.13%	0.015%
63	18.109	2487	2489	2493	rVB3	2266	2410	0.27%	0.029%
64	18.139	2493	2494	2497	rBV3	1494	1314	0.15%	0.016%
65	18.286	2515	2519	2522	rVB2	5323	6892	0.76%	0.084%

66	18.509	2552	2557	2560	rBV4	5827	9374	1.03%	0.114%
67	18.579	2567	2569	2573	rVB4	2269	1508	0.17%	0.018%
68	18.826	2609	2611	2612	rBV2	1771	1729	0.19%	0.021%
69	18.920	2625	2627	2630	rVB3	2487	1819	0.20%	0.022%
70	19.196	2671	2674	2677	rVB2	924	1264	0.14%	0.015%

71	19.231	2677	2680	2682	rBV3	1824	1996	0.22%	0.024%
72	19.255	2682	2684	2685	rBV	1211	1195	0.13%	0.015%
73	19.367	2701	2703	2705	rVB3	1995	1210	0.13%	0.015%
74	19.502	2724	2726	2729	rVB3	1644	1400	0.15%	0.017%
75	19.549	2730	2734	2737	rVV3	15062	20461	2.26%	0.249%

76	19.584	2737	2740	2743	rVB3	7134	8759	0.97%	0.106%
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 Title : SVOA CALIBRATION

77	19.654	2748	2752	2755	rBV2	5492	8174	0.90%	0.099%
78	19.825	2778	2781	2782	rBV2	1953	1833	0.20%	0.022%
79	19.907	2792	2795	2796	rBV2	2162	1975	0.22%	0.024%
80	20.013	2807	2813	2826	rVB	610728	905853	100.00%	11.010%
81	20.165	2837	2839	2842	rBV3	1698	2130	0.24%	0.026%
82	20.301	2860	2862	2864	rBV3	2507	2098	0.23%	0.025%
83	20.906	2963	2965	2970	rVB5	10653	14111	1.56%	0.172%
84	21.564	3072	3077	3084	rBV10	15219	38795	4.28%	0.472%
85	21.957	3137	3144	3153	rVB2	302833	510181	56.32%	6.201%
86	25.206	3687	3697	3713	rVB2	288079	826118	91.20%	10.041%
87	25.441	3727	3737	3753	rVB2	164023	534858	59.04%	6.501%
88	32.137	4875	4877	4878	rBV2	3620	2062	0.23%	0.025%

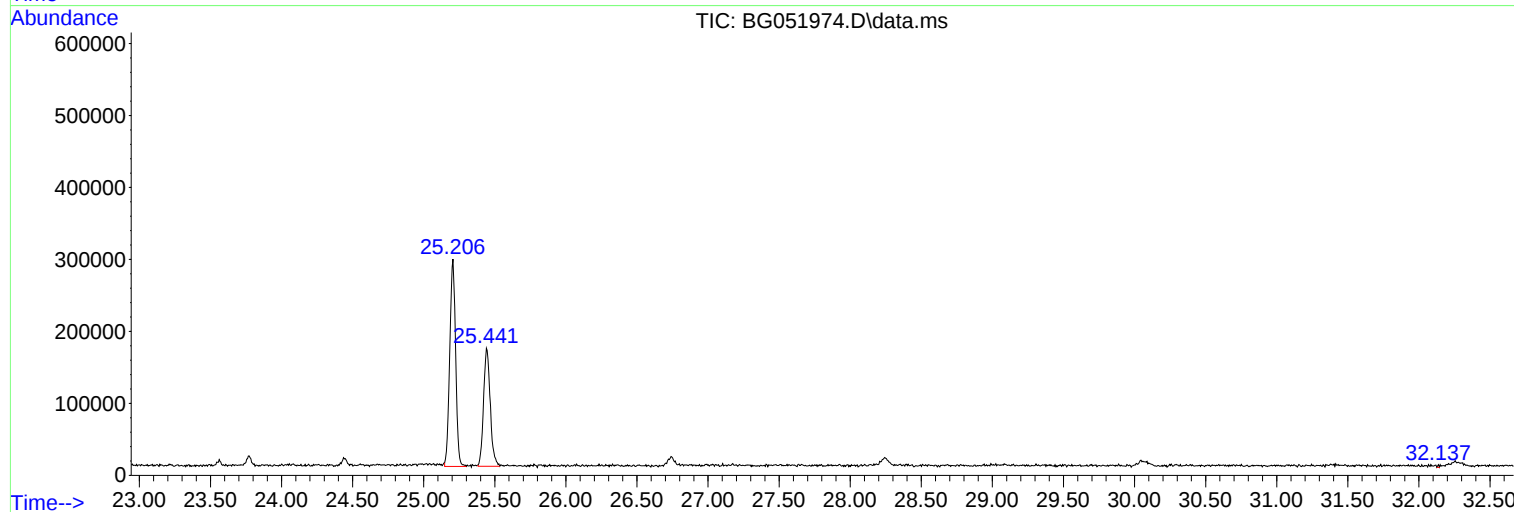
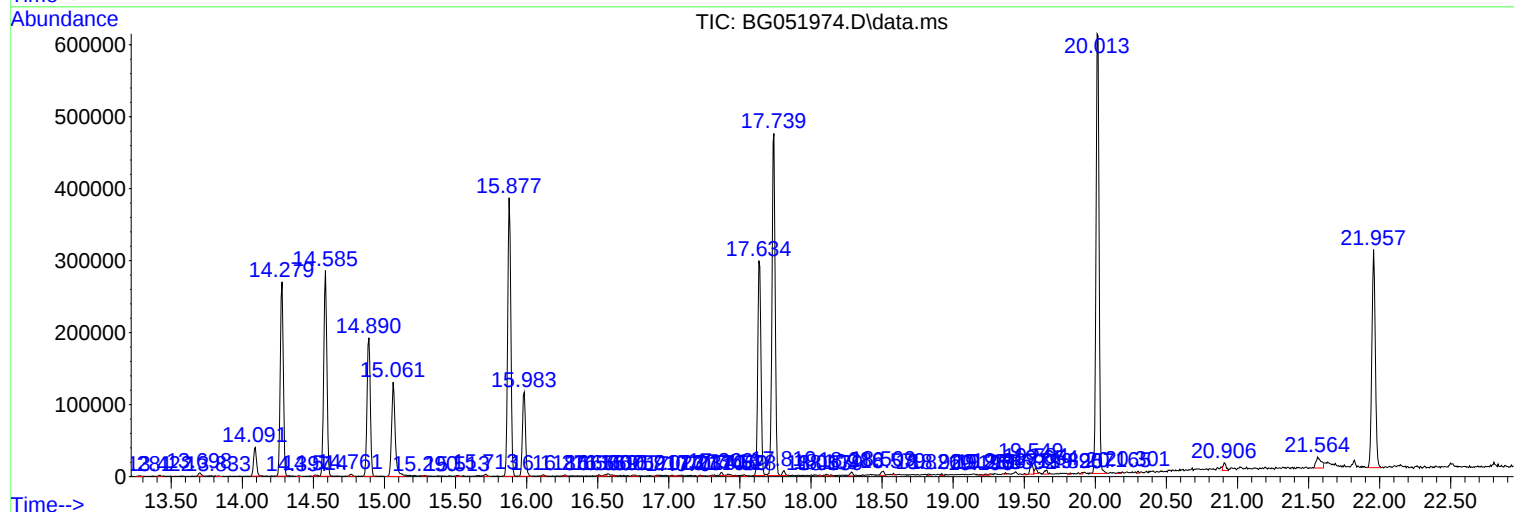
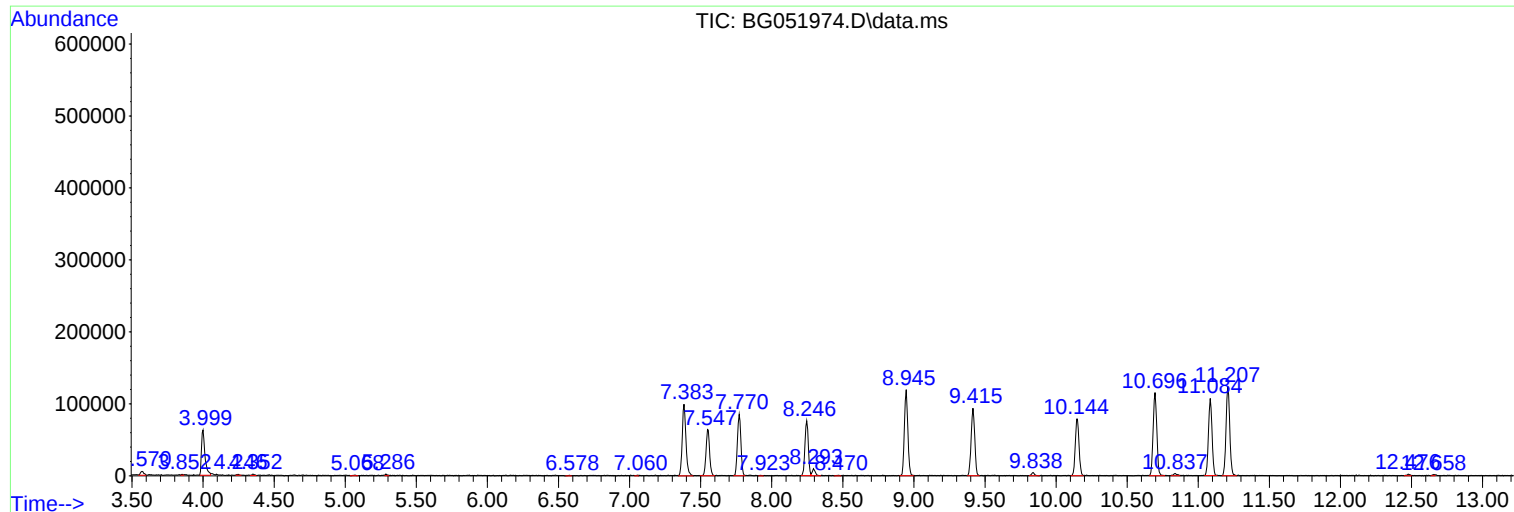
Sum of corrected areas: 8227713

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG010622.M
Quant Title : SVOA CALIBRATION

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



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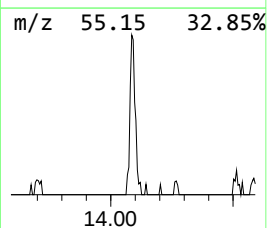
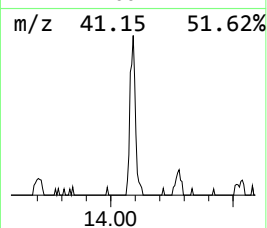
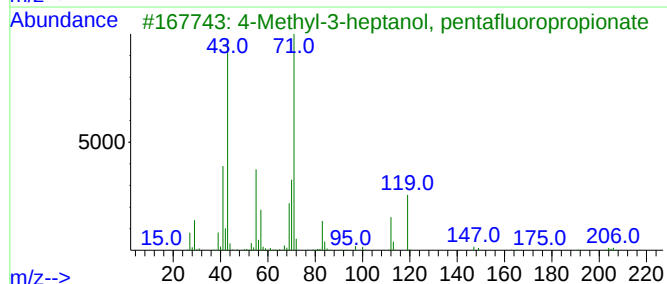
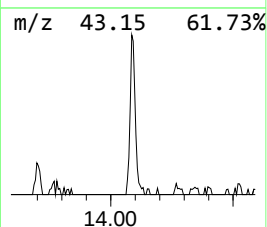
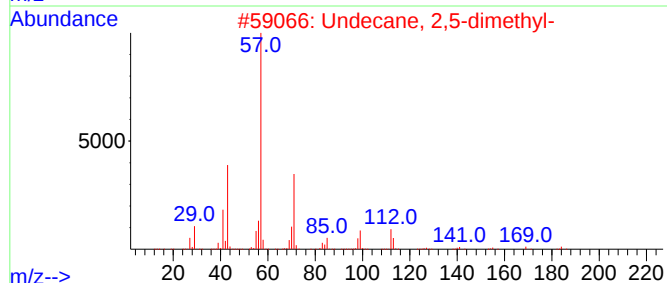
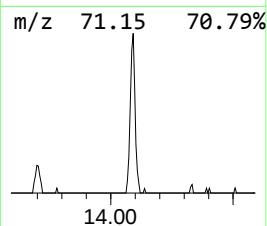
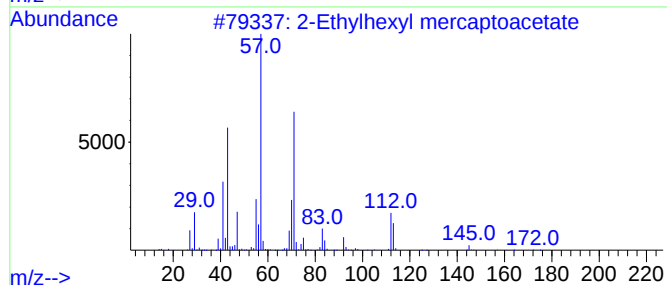
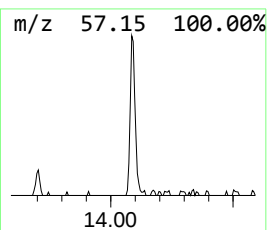
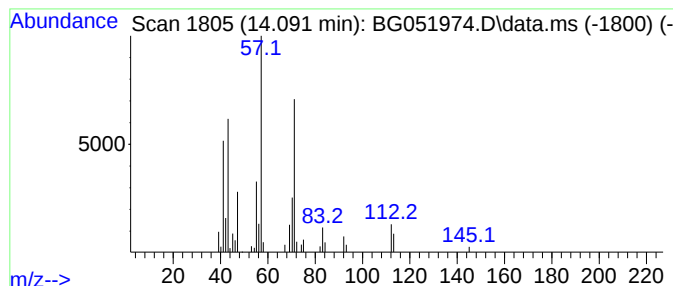
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG010622.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 2-Ethylhexyl mercaptoacetate Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.091	3.80 ng/ul	58209	Acenaphthene-d10	14.890

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Ethylhexyl mercaptoacetate	204	C10H20O2S	007659-86-1	78
2			Undecane, 2,5-dimethyl-	184	C13H28	017301-22-3	58
3			4-Methyl-3-heptanol, pentafluoro...	276	C11H17F5O2	1000365-51-7	53
4			4-Methyl-3-heptanol, trifluoroac...	226	C10H17F3O2	1000365-51-8	53
5			Octane, 3-ethyl-	142	C10H22	005881-17-4	50



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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
2-Ethylhexyl me...	14.091	3.8	ng/u1	58209	3	14.890	306183	20.0