

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101821\
 Data File : BP007541.D
 Acq On : 20 Oct 2021 22:05
 Operator : CG/JU
 Sample : M4246-13
 Misc :
 ALS Vial : 96 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 YAT55

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_P\METHODS\SFAM-EPA-BP101421.M
 Title : SVOA CALIBRATION

Signal : TIC: BP007541.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.375	12	15	24	rVB	47739	66120	1.14%	0.121%
2	3.657	59	63	67	rBV7	6543	12700	0.22%	0.023%
3	3.799	85	87	101	rBV	213346	329838	5.69%	0.601%
4	4.040	122	128	138	rVB	35718	53804	0.93%	0.098%
5	4.134	140	144	147	rBV2	13136	17950	0.31%	0.033%
6	4.228	156	160	165	rVB2	14181	18477	0.32%	0.034%
7	4.510	203	208	211	rVB7	5168	7358	0.13%	0.013%
8	5.051	296	300	305	rVB4	7485	10369	0.18%	0.019%
9	6.328	513	517	523	rVB2	11449	18555	0.32%	0.034%
10	7.110	643	650	659	rVB	226897	358292	6.18%	0.653%
11	7.281	670	679	687	rBV	676109	1092430	18.86%	1.991%
12	7.481	705	713	722	rBV	741055	1145495	19.77%	2.088%
13	7.587	729	731	738	rVB7	4618	7092	0.12%	0.013%
14	7.951	786	793	801	rVB	781692	1228842	21.21%	2.240%
15	8.204	830	836	845	rBV	5529	11136	0.19%	0.020%
16	8.516	883	889	892	rBV8	4238	6882	0.12%	0.013%
17	8.657	905	913	921	rBV	603302	963081	16.62%	1.755%
18	9.104	982	989	999	rBV	808744	1331736	22.99%	2.427%
19	9.222	1006	1009	1015	rVV8	5659	10013	0.17%	0.018%
20	9.287	1015	1020	1025	rVB9	4375	9140	0.16%	0.017%
21	9.575	1064	1069	1073	rVB5	8799	15563	0.27%	0.028%
22	9.834	1104	1113	1122	rBV	614784	1005356	17.35%	1.832%
23	10.375	1197	1205	1217	rBV	1000075	1670902	28.84%	3.046%
24	10.539	1225	1233	1241	rBV	251945	412780	7.13%	0.752%
25	10.757	1257	1270	1279	rBV	1056151	1769218	30.54%	3.225%
26	10.886	1283	1292	1302	rBV	800075	1360156	23.48%	2.479%
27	11.404	1370	1380	1383	rBV10	3881	8865	0.15%	0.016%
28	12.075	1486	1494	1499	rBV5	8407	16350	0.28%	0.030%
29	12.339	1534	1539	1544	rBV	22891	34230	0.59%	0.062%
30	12.433	1550	1555	1562	rVB6	13089	21752	0.38%	0.040%
31	12.963	1640	1645	1650	rBV4	10866	18385	0.32%	0.034%
32	13.210	1681	1687	1692	rBV3	19699	29820	0.51%	0.054%
33	13.433	1720	1725	1730	rBV	25970	37669	0.65%	0.069%
34	13.498	1730	1736	1743	rVV	161866	249202	4.30%	0.454%
35	13.557	1743	1746	1753	rVV8	6712	12250	0.21%	0.022%
36	13.628	1754	1758	1765	rVB5	10165	15798	0.27%	0.029%

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 Stop Thrs : 0 Peak Location: TOP

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Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP101421.M
 Title : SVOA CALIBRATION

37	13.792	1780	1786	1791	rVB9	9790	15734	0.27%	0.029%
38	13.998	1814	1821	1835	rBV	1902738	2709240	46.77%	4.938%
39	14.280	1858	1869	1878	rBV	2153863	3208417	55.38%	5.848%
40	14.586	1913	1921	1941	rBV	1658262	2560310	44.20%	4.667%
41	14.769	1945	1952	1964	rVV	209160	325386	5.62%	0.593%
42	15.004	1988	1992	1997	rVB8	9975	16575	0.29%	0.030%
43	15.274	2035	2038	2041	rVB4	5756	7317	0.13%	0.013%
44	15.404	2054	2060	2065	rBV6	7734	14983	0.26%	0.027%
45	15.580	2083	2090	2101	rVB	2920573	4175948	72.08%	7.611%
46	15.692	2102	2109	2118	rBV	701621	978089	16.88%	1.783%
47	15.822	2125	2131	2136	rBV	33970	48240	0.83%	0.088%
48	16.027	2163	2166	2169	rBV4	8520	12976	0.22%	0.024%
49	16.145	2181	2186	2194	rVB4	7777	15156	0.26%	0.028%
50	16.257	2199	2205	2211	rVV9	5296	12879	0.22%	0.023%
51	16.316	2211	2215	2218	rVV5	7404	10853	0.19%	0.020%
52	16.363	2221	2223	2230	rVB8	6348	11082	0.19%	0.020%
53	16.486	2239	2244	2249	rVV8	8612	13392	0.23%	0.024%
54	16.592	2259	2262	2266	rBV5	5376	9368	0.16%	0.017%
55	16.745	2284	2288	2292	rVB6	6924	8824	0.15%	0.016%
56	17.127	2346	2353	2356	rBV2	11379	19193	0.33%	0.035%
57	17.339	2382	2389	2395	rBV	2053560	2883290	49.77%	5.255%
58	17.439	2399	2406	2417	rBV2	3318948	4842568	83.59%	8.826%
59	17.716	2449	2453	2454	rBV3	4946	6661	0.11%	0.012%
60	17.739	2454	2457	2463	rVB8	9996	13825	0.24%	0.025%
61	18.021	2500	2505	2513	rBV	144886	180826	3.12%	0.330%
62	18.233	2534	2541	2550	rBV2	83665	137689	2.38%	0.251%
63	18.298	2550	2552	2558	rVB	9623	12134	0.21%	0.022%
64	19.121	2689	2692	2693	rBV2	14704	14759	0.25%	0.027%
65	19.151	2693	2697	2707	rVB	233303	277010	4.78%	0.505%
66	19.251	2710	2714	2716	rBV2	50795	62980	1.09%	0.115%
67	19.280	2716	2719	2722	rVV	121875	132100	2.28%	0.241%
68	19.315	2722	2725	2729	rVB	48403	58242	1.01%	0.106%
69	19.463	2747	2750	2755	rBV3	34196	41087	0.71%	0.075%
70	19.610	2773	2775	2778	rBV3	12229	13669	0.24%	0.025%
71	19.668	2779	2785	2792	rVV	4235706	5783414	99.83%	10.541%
72	21.145	3032	3036	3042	rBV4	113544	183792	3.17%	0.335%
73	21.421	3078	3083	3089	rBV	2490634	3392503	58.56%	6.183%
74	23.715	3466	3473	3482	rVB2	2883294	5793206	100.00%	10.559%
75	23.874	3493	3500	3511	rVB2	1622864	3483247	60.13%	6.349%

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Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP101421.M
Title : SVOA CALIBRATION

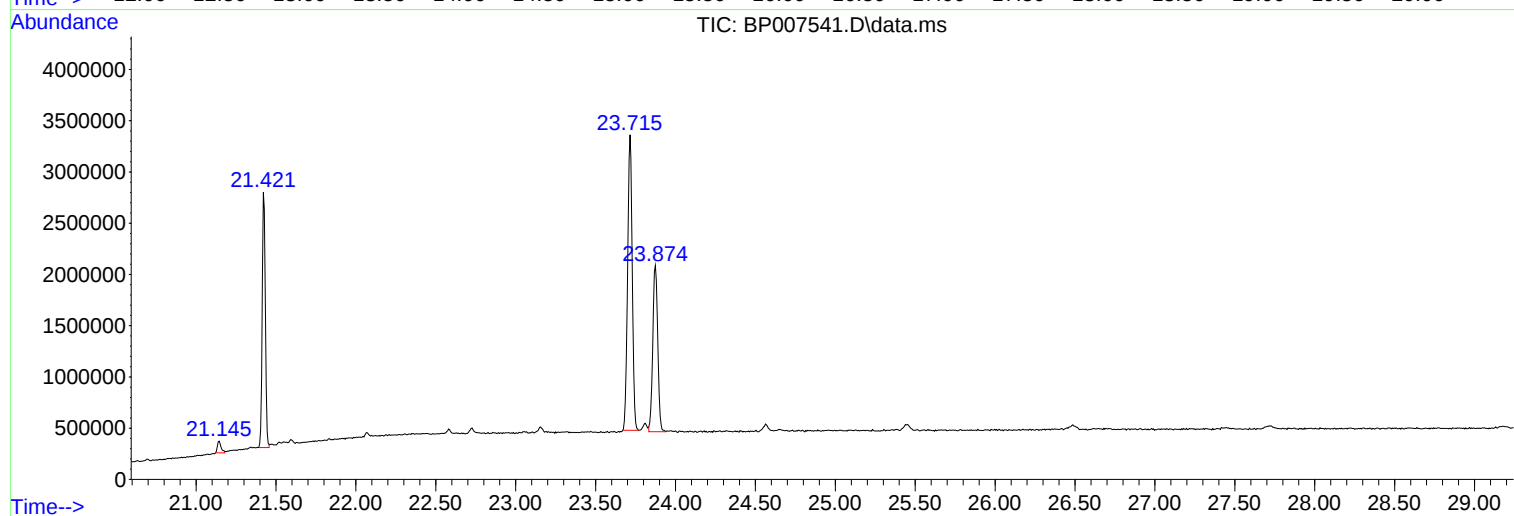
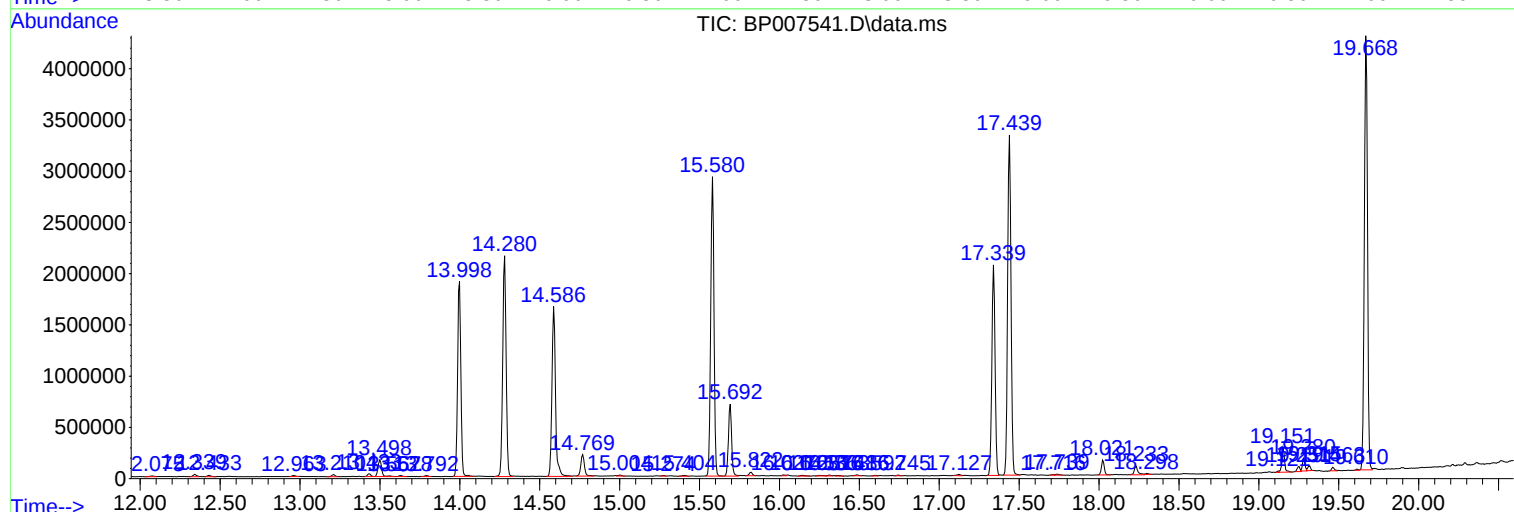
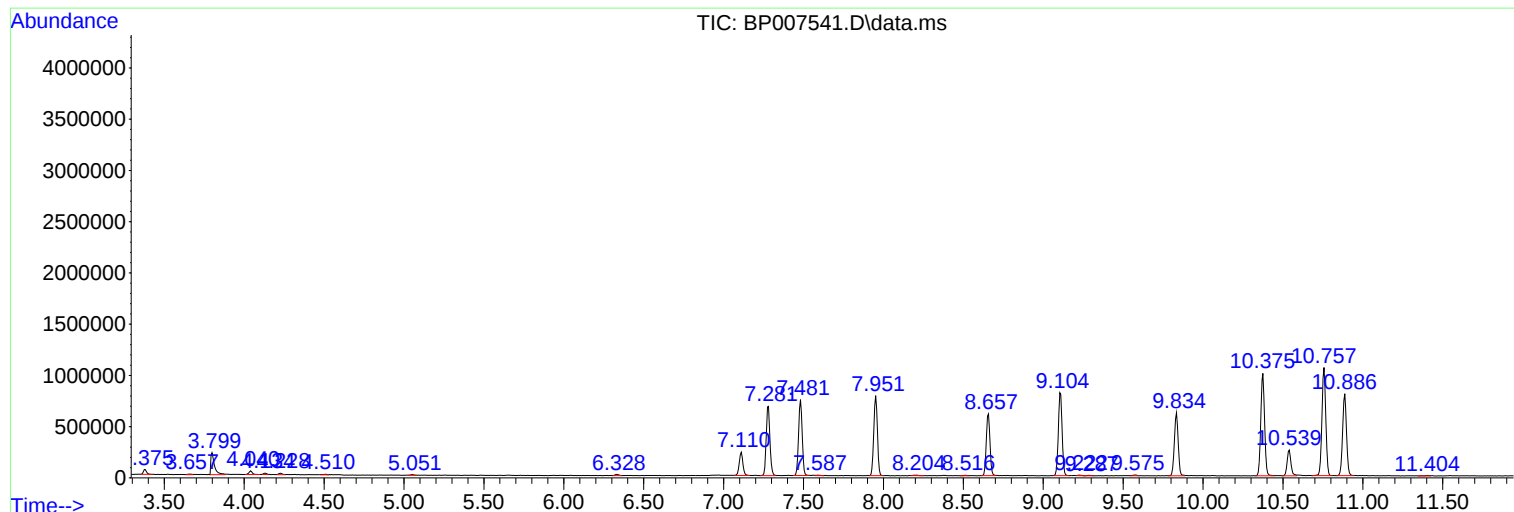
Sum of corrected areas: 54864570

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP101421.M
Quant Title : SVOA CALIBRATION

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



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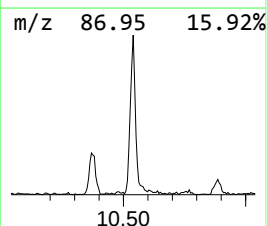
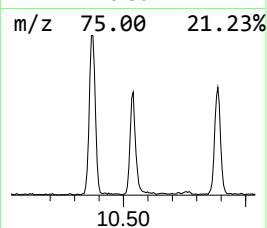
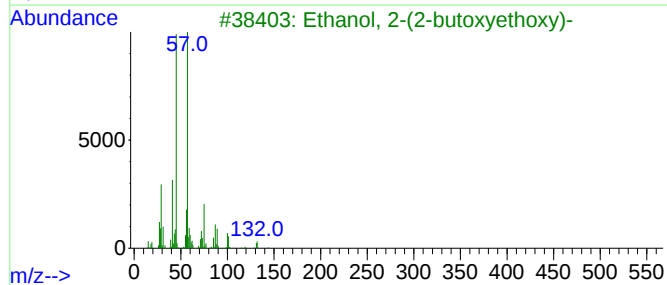
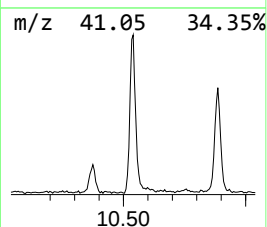
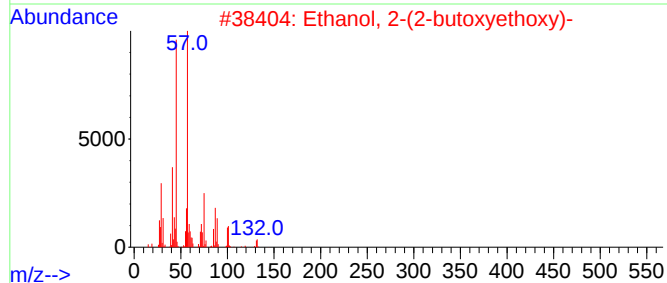
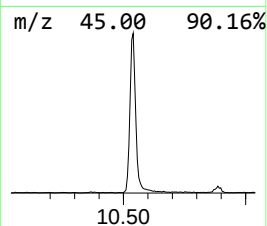
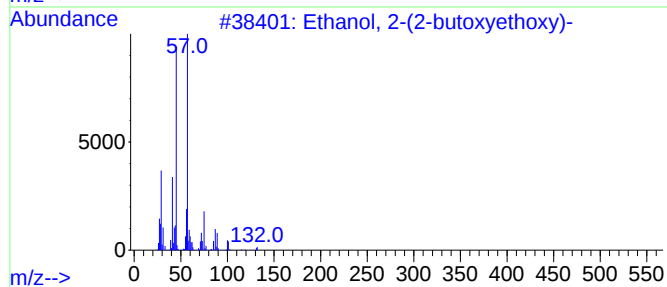
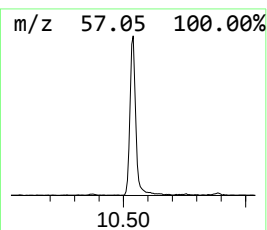
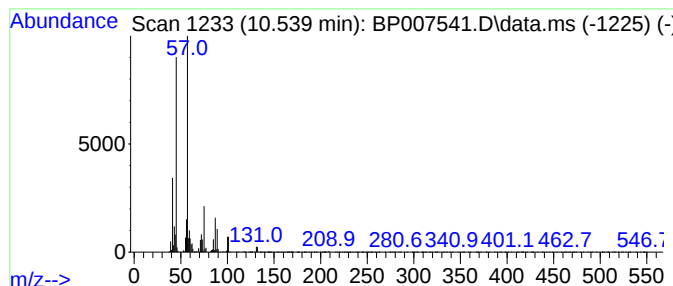
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP101421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.539	4.67 ng/ul	412780	Naphthalene-d8	10.757

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
2			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
3			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
4			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	83
5			Ethanol, 2-[2-(2-methylpropoxy)e...	162	C8H18O3	018912-80-6	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
Ethanol, 2-(2-b...	10.539	4.7	ng/u1	412780	2	10.757	1769220	20.0