

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
 Data File : BP012352.D
 Acq On : 27 Oct 2022 20:40
 Operator : CG/JU
 Sample : N5015-12
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C1BE8

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-BP102522.M
 Title : SVOA CALIBRATION

Signal : TIC: BP012352.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.252	41	45	55	rVB	82954	124185	1.15%	0.116%
2	3.540	92	94	104	rVB	17452	35721	0.33%	0.033%
3	3.681	116	118	132	rBV	183625	337219	3.13%	0.316%
4	3.775	132	134	138	rVB4	13469	14921	0.14%	0.014%
5	3.887	148	153	161	rVB	68923	107411	1.00%	0.101%
6	3.987	166	170	174	rVB	20446	25612	0.24%	0.024%
7	4.364	230	234	240	rVB2	19190	30292	0.28%	0.028%
8	4.558	262	267	270	rVB7	7091	11792	0.11%	0.011%
9	4.905	322	326	338	rVB2	50213	105060	0.97%	0.099%
10	6.981	672	679	698	rBV	1346876	2377633	22.05%	2.229%
11	7.146	698	707	725	rVB	1208204	1957984	18.16%	1.836%
12	7.340	733	740	750	rBV	1398738	2275963	21.10%	2.134%
13	7.410	750	752	757	rVB3	9403	13251	0.12%	0.012%
14	7.634	785	790	796	rVB3	9118	15057	0.14%	0.014%
15	7.805	812	819	828	rBV	1110386	1807728	16.76%	1.695%
16	8.522	934	941	956	rBV	1717079	2852360	26.45%	2.674%
17	8.640	956	961	969	rVB	49986	93468	0.87%	0.088%
18	8.975	1011	1018	1035	rBV	1388161	2309559	21.42%	2.166%
19	9.363	1078	1084	1089	rBV3	15781	26860	0.25%	0.025%
20	9.699	1132	1141	1153	rBV	1100434	1900963	17.63%	1.782%
21	10.057	1199	1202	1212	rVB	4995	11109	0.10%	0.010%
22	10.234	1225	1232	1253	rBV	1754737	3074644	28.51%	2.883%
23	10.399	1253	1260	1288	rVV	368777	905013	8.39%	0.849%
24	10.610	1288	1296	1311	rVB	1657429	2869321	26.61%	2.690%
25	10.757	1314	1321	1341	rBV	1562436	2830120	26.24%	2.654%
26	11.451	1431	1439	1444	rBV	7403	19877	0.18%	0.019%
27	12.204	1561	1567	1576	rVB2	31867	58064	0.54%	0.054%
28	12.804	1664	1669	1674	rBV8	5921	11087	0.10%	0.010%
29	13.063	1708	1713	1719	rVB5	9627	13503	0.13%	0.013%
30	13.181	1728	1733	1738	rBV6	7213	12277	0.11%	0.012%
31	13.269	1743	1748	1751	rVV6	10112	17087	0.16%	0.016%
32	13.322	1753	1757	1759	rVV5	14095	23294	0.22%	0.022%
33	13.345	1759	1761	1769	rVB	25291	38786	0.36%	0.036%
34	13.422	1769	1774	1781	rBV3	10777	18630	0.17%	0.017%
35	13.787	1830	1836	1842	rBV2	21663	42160	0.39%	0.040%
36	13.875	1844	1851	1865	rVV	3618455	5028478	46.63%	4.715%

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Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
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37	13.969	1865	1867	1875	rVB9	8604	15639	0.15%	0.015%
38	14.151	1885	1898	1915	rBV	4165811	6134977	56.89%	5.752%
39	14.345	1925	1931	1937	rVB	13070	23960	0.22%	0.022%
40	14.457	1941	1950	1963	rBV	2956739	4223915	39.17%	3.960%
41	14.628	1975	1979	1981	rBV3	9185	11911	0.11%	0.011%
42	14.675	1981	1987	2009	rVV	1316056	2280472	21.15%	2.138%
43	14.887	2018	2023	2032	rVB	133589	218382	2.02%	0.205%
44	14.963	2035	2036	2046	rVB9	8325	14152	0.13%	0.013%
45	15.298	2088	2093	2098	rVB2	13038	22066	0.20%	0.021%
46	15.369	2101	2105	2109	rVB7	8552	11182	0.10%	0.010%
47	15.457	2113	2120	2133	rBV	5304115	7868476	72.96%	7.378%
48	15.587	2136	2142	2156	rBV	1431282	2023889	18.77%	1.898%
49	15.692	2156	2160	2167	rVB	29235	49264	0.46%	0.046%
50	15.916	2193	2198	2202	rBV5	18938	31699	0.29%	0.030%
51	16.028	2212	2217	2223	rBV10	8986	15470	0.14%	0.015%
52	16.498	2291	2297	2300	rBV8	6886	16549	0.15%	0.016%
53	16.645	2319	2322	2327	rVB7	14812	23265	0.22%	0.022%
54	16.722	2332	2335	2338	rVV3	13102	17624	0.16%	0.017%
55	16.757	2338	2341	2346	rVB	25227	33055	0.31%	0.031%
56	17.004	2379	2383	2385	rBV4	7808	11908	0.11%	0.011%
57	17.133	2401	2405	2410	rVB8	8854	15472	0.14%	0.015%
58	17.216	2413	2419	2425	rBV	3553619	5269927	48.87%	4.941%
59	17.322	2430	2437	2446	rVV2	5828583	8661823	80.32%	8.122%
60	17.422	2450	2454	2461	rVB2	103084	148754	1.38%	0.139%
61	17.886	2529	2533	2538	rVB	83499	99066	0.92%	0.093%
62	17.986	2547	2550	2554	rBV5	20709	34678	0.32%	0.033%
63	18.104	2566	2570	2580	rBV	311555	490864	4.55%	0.460%
64	19.133	2741	2745	2751	rBV	94202	143659	1.33%	0.135%
65	19.204	2753	2757	2759	rVV	97893	124041	1.15%	0.116%
66	19.233	2759	2762	2768	rVB	100761	145773	1.35%	0.137%
67	19.339	2776	2780	2788	rBV	228862	359889	3.34%	0.337%
68	19.557	2812	2817	2831	rBV	7638653	10784521	100.00%	10.112%
69	20.704	3009	3012	3017	rBV3	93272	151849	1.41%	0.142%
70	21.022	3062	3066	3076	rBV2	314083	683011	6.33%	0.640%
71	21.263	3103	3107	3110	rBV	340465	414158	3.84%	0.388%
72	21.316	3111	3116	3127	rVB	6399249	8227112	76.29%	7.714%
73	22.110	3248	3251	3259	rVB2	361374	531105	4.92%	0.498%
74	22.957	3391	3395	3402	rVB	317504	500712	4.64%	0.469%
75	23.557	3490	3497	3512	rVB2	4856263	10156766	94.18%	9.523%
76	23.710	3517	3523	3533	rVB2	2593257	5264387	48.81%	4.936%

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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_P\Methods\SFAM-EPA-BP102522.M
Title : SVOA CALIBRATION

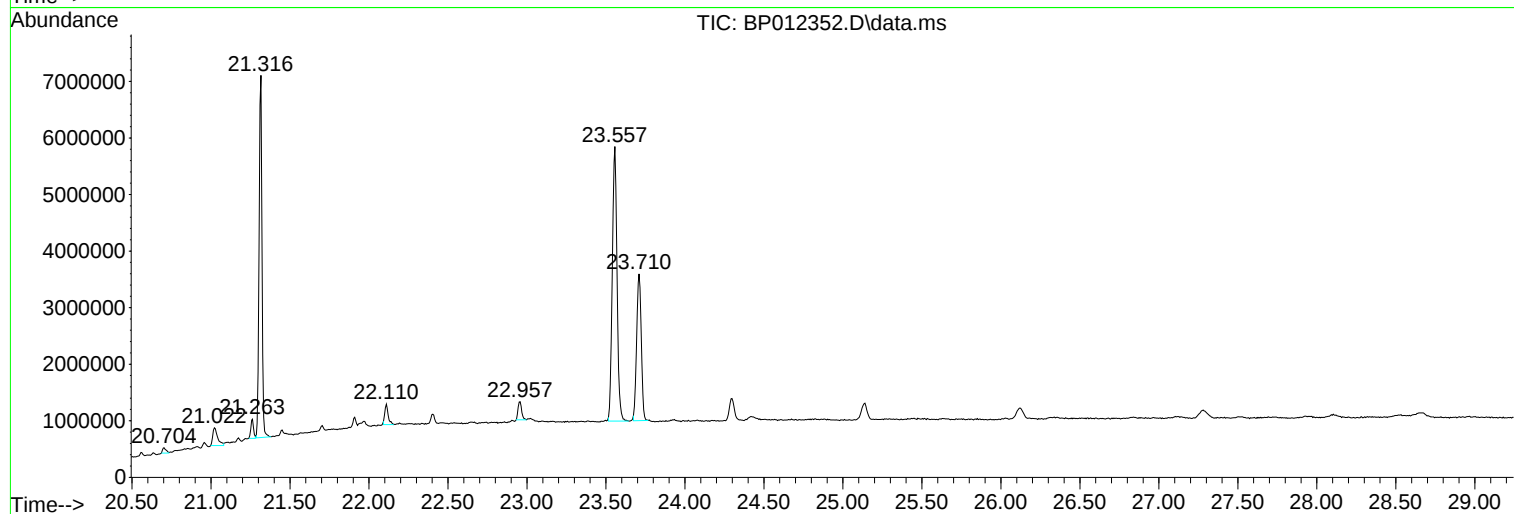
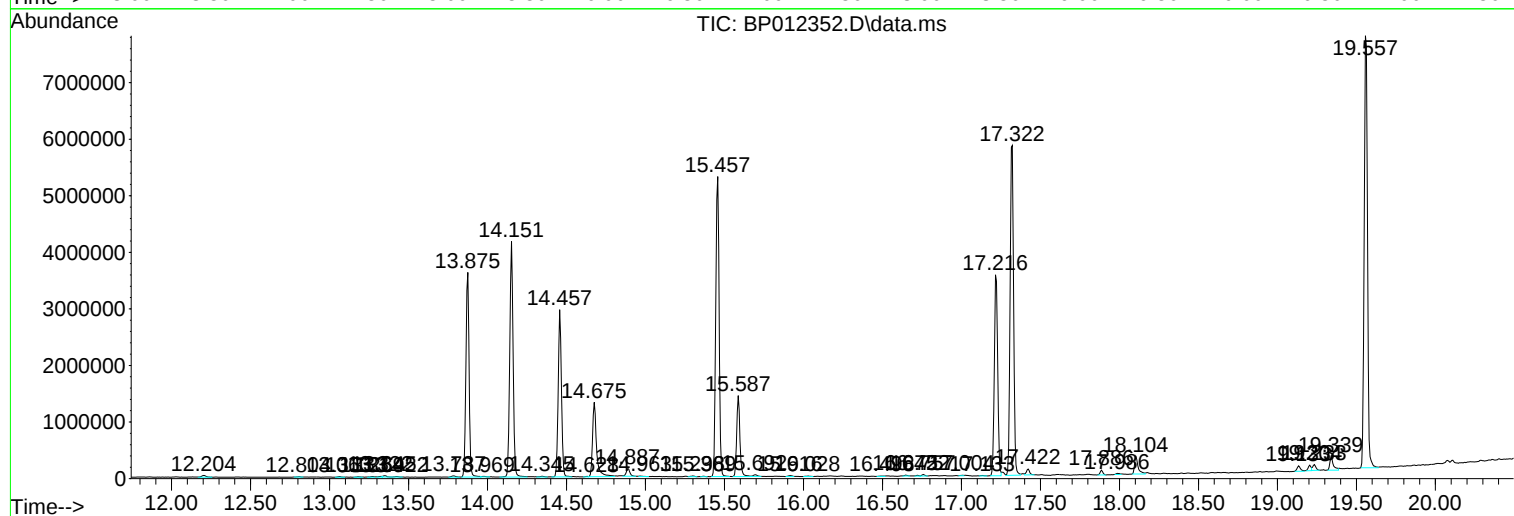
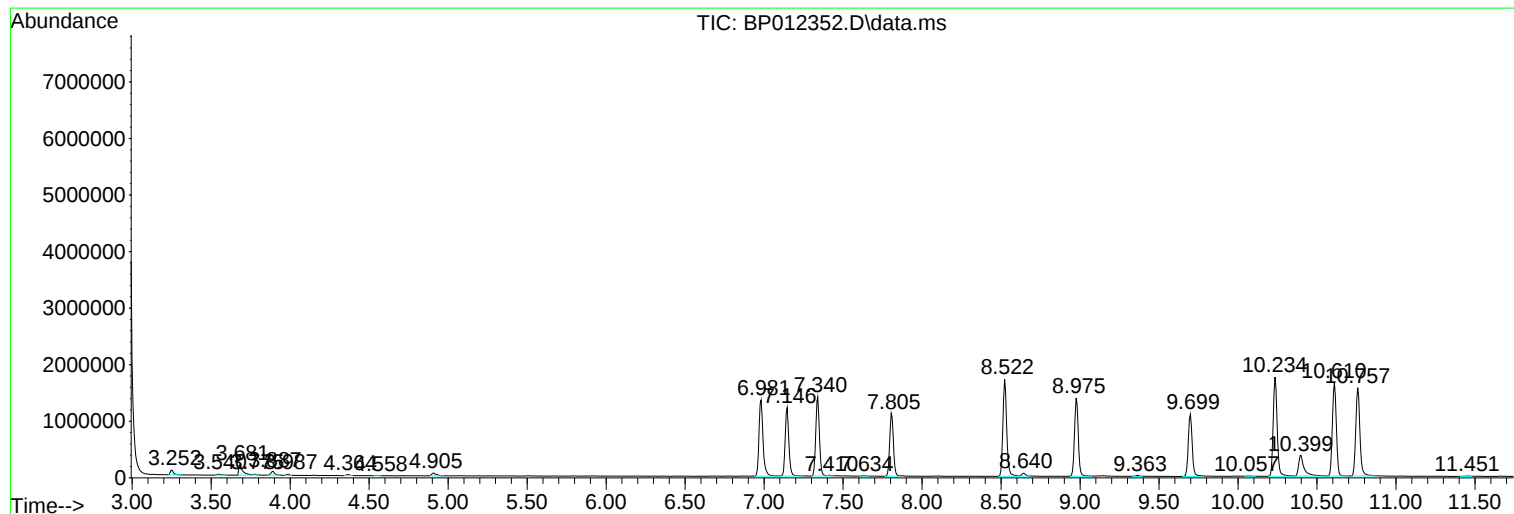
Sum of corrected areas: 106651911

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Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

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



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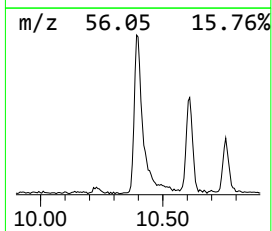
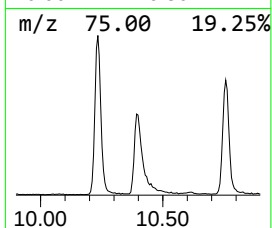
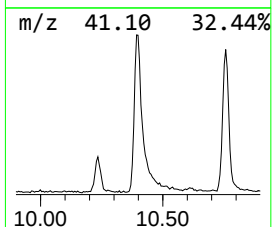
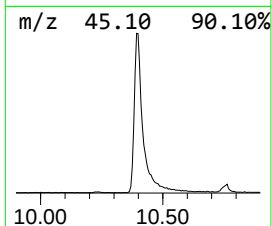
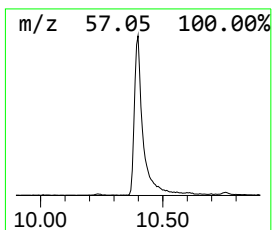
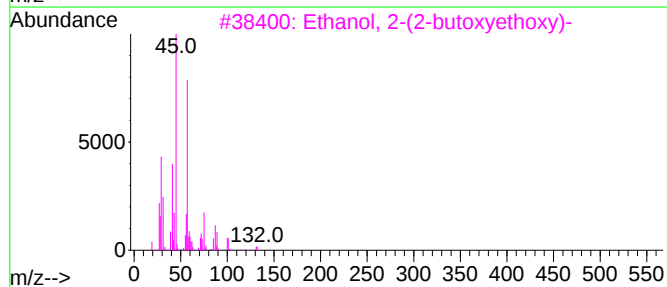
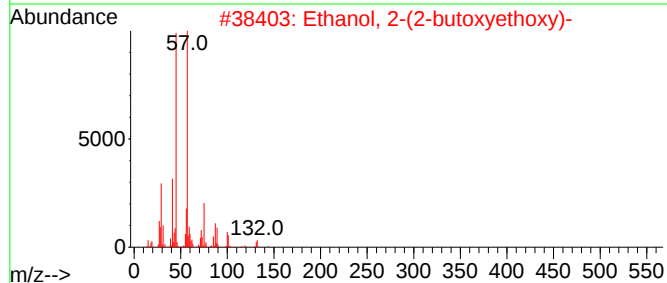
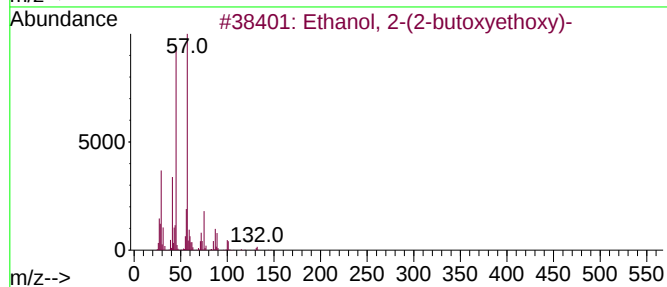
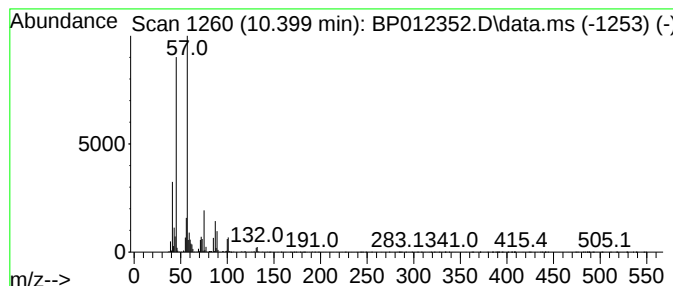
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST0.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.
10.399	6.31 ng/ul	905013	Naphthalene-d8	10.610

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, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	83
5			Ethylene glycol monoisobutyl ether	118	C6H14O2	004439-24-1	64



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TIC Library : C:\DATABASE\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Ethanol, 2-(2-b...	10.399	6.3	ng/u1	905013	2	10.610	2869320	20.0