

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110222\
 Data File : BM037316.D
 Acq On : 02 Nov 2022 15:02
 Operator : CG/JU
 Sample : N5301-09
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBWU6

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_M\Methods\SFAM-EPA-BM110122.M
 Title : SVOA CALIBRATION

Signal : TIC: BM037316.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.240	22	26	29	rBV	22378	29926	0.35%	0.037%
2	3.275	29	32	37	rVB	64426	73672	0.87%	0.092%
3	3.699	101	104	117	rBV	460161	675235	7.94%	0.840%
4	3.922	137	142	147	rVB2	25121	38457	0.45%	0.048%
5	4.022	154	159	164	rVB	29142	40351	0.47%	0.050%
6	4.969	315	320	326	rBV	18051	29482	0.35%	0.037%
7	5.563	414	421	426	rVB8	7867	16130	0.19%	0.020%
8	6.969	654	660	663	rBV4	8655	14950	0.18%	0.019%
9	7.028	664	670	683	rVV	293037	498315	5.86%	0.620%
10	7.116	683	685	689	rVB3	7447	10854	0.13%	0.013%
11	7.193	692	698	712	rBV	1098582	1692665	19.91%	2.105%
12	7.387	724	731	741	rBV	1004279	1579462	18.58%	1.964%
13	7.463	741	744	746	rVV4	8662	10754	0.13%	0.013%
14	7.640	768	774	784	rBV	88160	158295	1.86%	0.197%
15	7.857	804	811	819	rBV	1156085	1814565	21.35%	2.257%
16	7.945	824	826	833	rVB8	5206	11236	0.13%	0.014%
17	8.575	926	933	944	rBV	846640	1425945	16.77%	1.773%
18	8.669	946	949	954	rVB4	11699	19885	0.23%	0.025%
19	9.022	1002	1009	1022	rBV	1242935	2047037	24.08%	2.546%
20	9.198	1034	1039	1044	rVB10	5801	12478	0.15%	0.016%
21	9.369	1064	1068	1071	rBV6	5432	9711	0.11%	0.012%
22	9.410	1071	1075	1082	rVV	27547	50435	0.59%	0.063%
23	9.745	1125	1132	1148	rBV	940241	1539083	18.10%	1.914%
24	10.186	1202	1207	1213	rBV2	16732	27827	0.33%	0.035%
25	10.281	1216	1223	1243	rBV	1540081	2528986	29.75%	3.145%
26	10.439	1244	1250	1263	rBV	431534	794636	9.35%	0.988%
27	10.657	1280	1287	1295	rBV	1697321	2757265	32.43%	3.429%
28	10.804	1305	1312	1328	rBV	1358663	2311723	27.19%	2.875%
29	10.904	1328	1329	1334	rVB5	7968	8667	0.10%	0.011%
30	11.275	1389	1392	1398	rBV7	6009	12365	0.15%	0.015%
31	11.333	1400	1402	1407	rVB6	6921	9136	0.11%	0.011%
32	11.510	1423	1432	1434	rBV10	4866	11807	0.14%	0.015%
33	11.792	1475	1480	1486	rBV6	9420	15207	0.18%	0.019%
34	11.904	1491	1499	1506	rVB	34217	60789	0.72%	0.076%
35	12.245	1553	1557	1563	rBV2	25746	42450	0.50%	0.053%
36	12.422	1583	1587	1591	rBV7	6602	9804	0.12%	0.012%

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Min Area: 0.1 % of largest Peak

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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_M\Methods\SFAM-EPA-BM110122.M
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37	12.698	1630	1634	1640	rVB7	9305	16383	0.19%	0.020%
38	12.786	1646	1649	1653	rBV4	12079	19946	0.23%	0.025%
39	12.851	1653	1660	1664	rVB9	7297	18454	0.22%	0.023%
40	13.098	1698	1702	1709	rVB8	10873	22409	0.26%	0.028%
41	13.227	1719	1724	1731	rBV8	6779	13054	0.15%	0.016%
42	13.386	1746	1751	1756	rBV	53073	82561	0.97%	0.103%
43	13.504	1765	1771	1782	rVV	123812	208050	2.45%	0.259%
44	13.616	1786	1790	1797	rVV7	15916	26712	0.31%	0.033%
45	13.910	1832	1840	1847	rBV	3262862	4375014	51.46%	5.441%
46	14.110	1869	1874	1880	rBV4	28327	56282	0.66%	0.070%
47	14.186	1880	1887	1895	rBV	3824808	5394892	63.46%	6.710%
48	14.374	1915	1919	1922	rBV6	5235	8579	0.10%	0.011%
49	14.492	1932	1939	1947	rBV	2647816	3796991	44.67%	4.722%
50	14.716	1971	1977	1991	rBV	174602	391409	4.60%	0.487%
51	15.145	2047	2050	2054	rVB6	9844	13185	0.16%	0.016%
52	15.239	2061	2066	2069	rBV3	12759	20544	0.24%	0.026%
53	15.280	2069	2073	2075	rVV2	14308	23359	0.27%	0.029%
54	15.316	2075	2079	2085	rVB	47596	75304	0.89%	0.094%
55	15.386	2087	2091	2096	rBV	42494	66037	0.78%	0.082%
56	15.480	2101	2107	2115	rVV	4986312	6999009	82.33%	8.705%
57	15.610	2123	2129	2142	rBV	1412573	1931160	22.72%	2.402%
58	15.721	2144	2148	2152	rVB2	25393	40246	0.47%	0.050%
59	16.045	2201	2203	2209	rVB7	8297	14296	0.17%	0.018%
60	16.192	2225	2228	2230	rBV4	8531	9312	0.11%	0.012%
61	16.263	2238	2240	2244	rVB5	12271	12836	0.15%	0.016%
62	16.368	2256	2258	2264	rVB8	5795	8585	0.10%	0.011%
63	16.439	2266	2270	2274	rBV3	27711	38246	0.45%	0.048%
64	16.686	2308	2312	2315	rBV5	13415	21306	0.25%	0.026%
65	16.986	2360	2363	2366	rBV4	9196	10481	0.12%	0.013%
66	17.139	2387	2389	2395	rVB7	10337	18198	0.21%	0.023%
67	17.233	2399	2405	2414	rVV	3155571	4294808	50.52%	5.342%
68	17.333	2415	2422	2432	rBV	5313583	7594070	89.33%	9.445%
69	17.898	2514	2518	2524	rVB	99570	124446	1.46%	0.155%
70	18.004	2533	2536	2541	rBV7	13493	23028	0.27%	0.029%
71	18.121	2552	2556	2566	rBV3	129888	214983	2.53%	0.267%
72	19.068	2714	2717	2722	rVB	203433	206283	2.43%	0.257%
73	19.186	2733	2737	2739	rBV2	63750	95827	1.13%	0.119%
74	19.256	2745	2749	2752	rBV	66973	98763	1.16%	0.123%
75	19.398	2769	2773	2782	rBV2	127326	223058	2.62%	0.277%
76	19.621	2805	2811	2817	rBV	6527960	8501024	100.00%	10.573%

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Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

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77	21.403	3109	3114	3123	rBV	3508503	4237859	49.85%	5.271%
78	23.603	3481	3488	3502	rVB	3592802	6936006	81.59%	8.626%
79	23.750	3507	3513	3524	rVB2	1886365	3731530	43.90%	4.641%

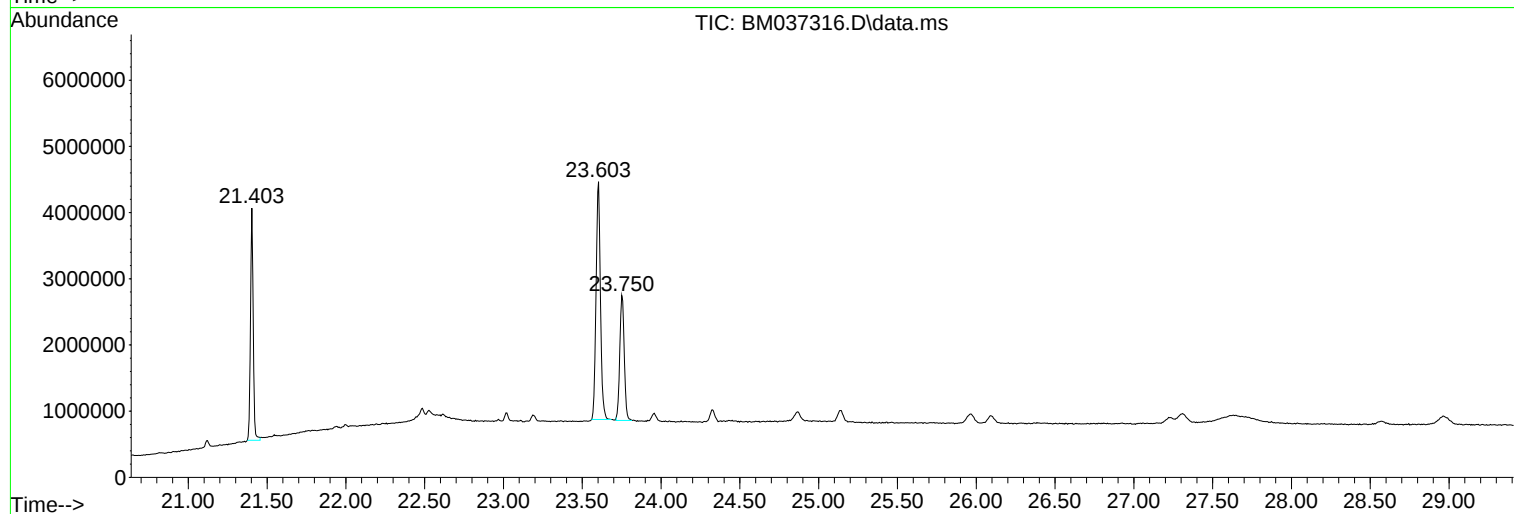
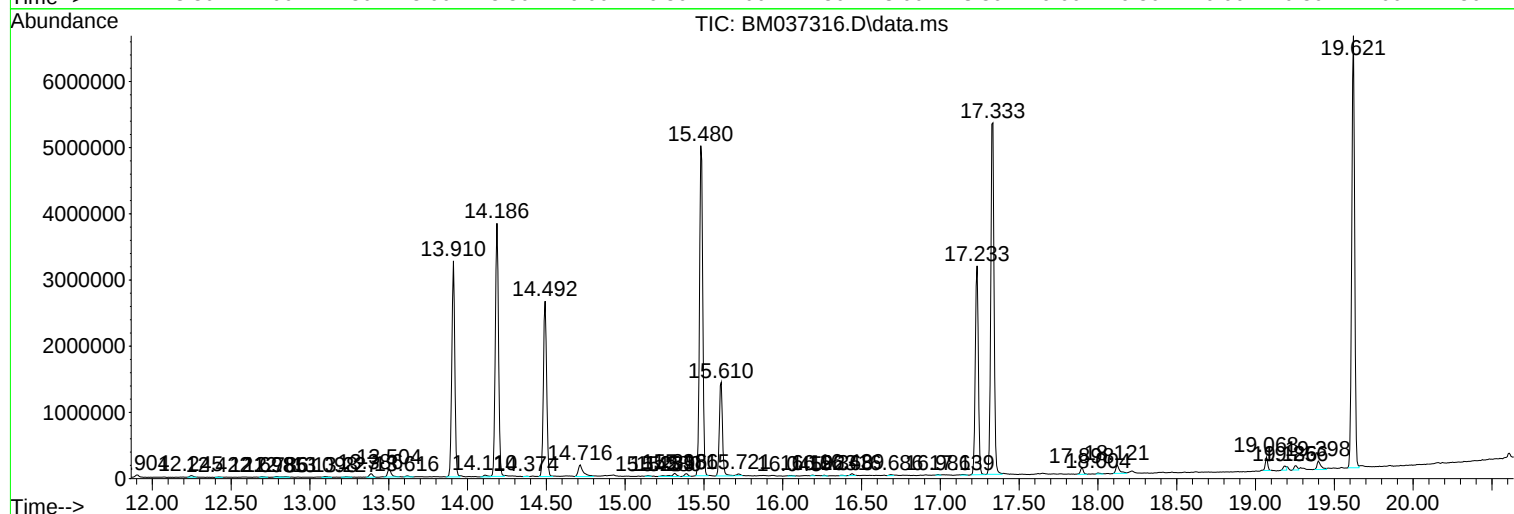
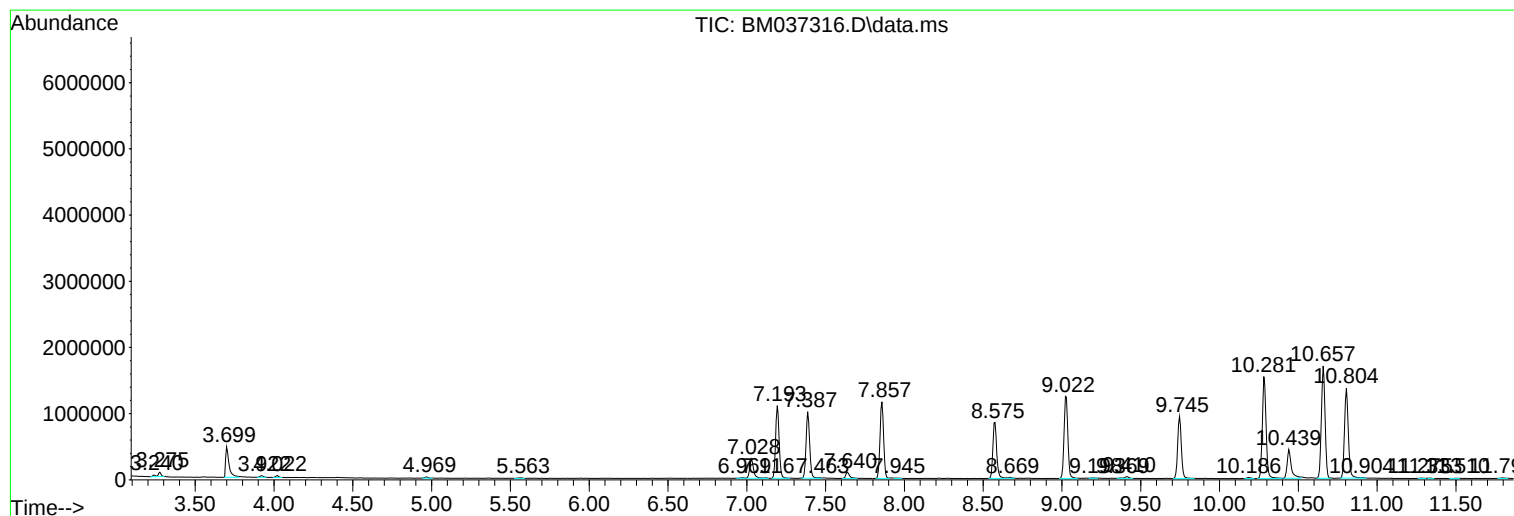
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM110122.M
Quant Title : SVOA CALIBRATION

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



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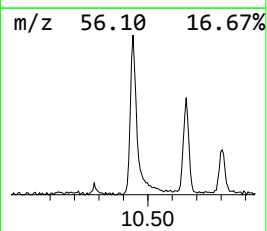
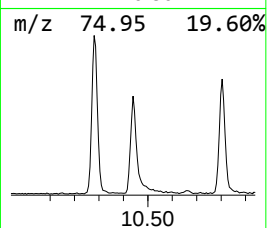
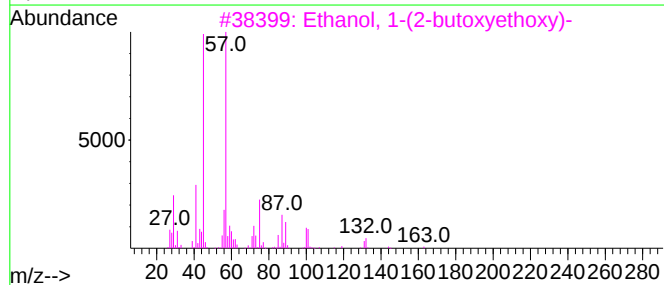
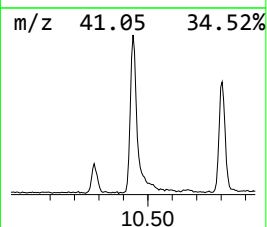
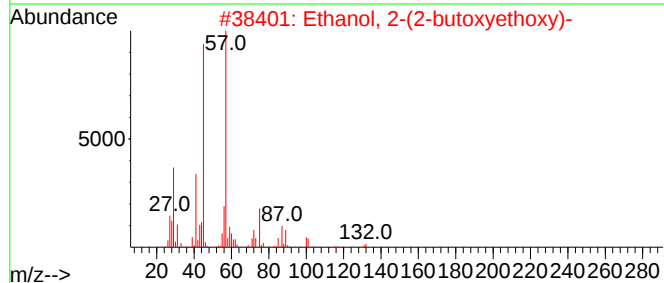
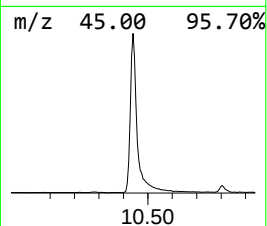
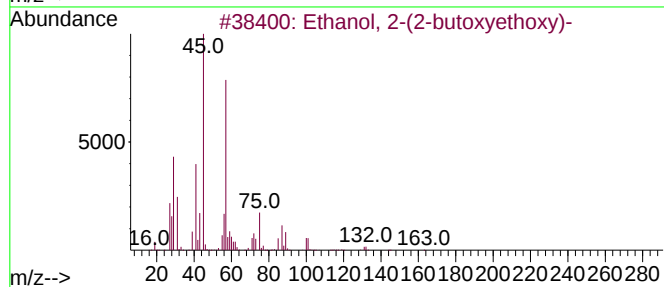
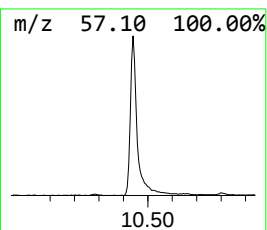
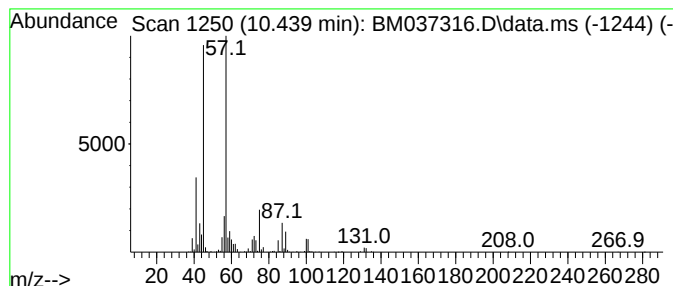
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM110122.M
Quant Title : SVOA CALIBRATION

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.
10.439	5.76 ng/ul	794636	Naphthalene-d8	10.657

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, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	83
4			Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	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...	10.439	5.8	ng/u1	794636	2	10.657	2757270	20.0