

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM083021\
 Data File : BM031956.D
 Acq On : 30 Aug 2021 18:29
 Operator : CG/JU
 Sample : M3365-06
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBK89

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-BM082821.M
 Title : SVOA CALIBRATION

Signal : TIC: BM031956.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.134	5	8	16	rVB	9280	15383	0.87%	0.087%
2	3.534	72	76	89	rBV	65791	135564	7.64%	0.767%
3	3.728	103	109	114	rVB	11369	19259	1.08%	0.109%
4	3.816	121	124	129	rVB4	6243	9269	0.52%	0.052%
5	4.028	157	160	165	rBV4	6035	9290	0.52%	0.053%
6	4.422	223	227	236	rBV4	3673	7757	0.44%	0.044%
7	4.693	270	273	280	rBV3	5751	8165	0.46%	0.046%
8	5.551	416	419	424	rBV5	1901	2848	0.16%	0.016%
9	6.704	608	615	628	rBV	166321	273812	15.42%	1.549%
10	6.869	637	643	654	rBV	120304	199759	11.25%	1.130%
11	6.940	654	655	662	rVB6	1886	2289	0.13%	0.013%
12	7.051	668	674	684	rBV	168583	283929	15.99%	1.606%
13	7.128	684	687	692	rVB3	2689	4263	0.24%	0.024%
14	7.510	745	752	761	rBV	337114	537583	30.28%	3.041%
15	7.581	763	764	771	rVB3	2650	2994	0.17%	0.017%
16	8.222	867	873	884	rBV	200580	337421	19.01%	1.909%
17	8.334	889	892	898	rVB3	5346	8206	0.46%	0.046%
18	8.657	941	947	957	rBV	165472	281784	15.87%	1.594%
19	8.792	967	970	974	rBV5	1863	2848	0.16%	0.016%
20	9.051	1009	1014	1017	rBV4	1687	2445	0.14%	0.014%
21	9.369	1063	1068	1080	rVB	144209	239563	13.49%	1.355%
22	9.904	1151	1159	1173	rBV	209178	383114	21.58%	2.167%
23	9.998	1173	1175	1180	rVB3	1550	1801	0.10%	0.010%
24	10.069	1181	1187	1205	rBV	135651	318972	17.97%	1.805%
25	10.263	1213	1220	1229	rVB	471199	799381	45.03%	4.522%
26	10.422	1240	1247	1262	rBV	198251	374459	21.09%	2.118%
27	11.874	1487	1494	1498	rBV4	3191	6138	0.35%	0.035%
28	12.751	1638	1643	1647	rBV5	2819	4160	0.23%	0.024%
29	12.974	1680	1681	1688	rVB4	1473	1823	0.10%	0.010%
30	13.051	1688	1694	1701	rVV3	3649	6628	0.37%	0.037%
31	13.574	1777	1783	1794	rBV	476070	657020	37.01%	3.717%
32	13.839	1821	1828	1836	rBV	504133	740464	41.71%	4.189%
33	14.145	1874	1880	1895	rVB	816844	1219137	68.67%	6.897%
34	14.339	1905	1913	1917	rBV6	2525	4280	0.24%	0.024%
35	14.398	1917	1923	1940	rBV	97016	244535	13.77%	1.383%
36	14.610	1955	1959	1963	rBV3	5963	8523	0.48%	0.048%

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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-BM082821.M
 Title : SVOA CALIBRATION

37	14.968	2017	2020	2023	rVB4	2169	2309	0.13%	0.013%
38	15.004	2023	2026	2030	rBV4	1569	1864	0.10%	0.011%
39	15.151	2045	2051	2062	rBV	716884	996580	56.14%	5.638%
40	15.298	2070	2076	2085	rBV	165657	246304	13.87%	1.393%
41	15.392	2088	2092	2096	rVB	6895	9284	0.52%	0.053%
42	15.621	2126	2131	2137	rVB7	5638	7724	0.44%	0.044%
43	15.898	2177	2178	2183	rVB3	1662	1869	0.11%	0.011%
44	16.068	2203	2207	2214	rBV6	1913	4205	0.24%	0.024%
45	16.468	2272	2275	2278	rVB2	1808	1784	0.10%	0.010%
46	16.686	2308	2312	2316	rBV4	4320	6612	0.37%	0.037%
47	16.904	2343	2349	2360	rVV	1077236	1529652	86.16%	8.654%
48	17.004	2360	2366	2379	rVV2	729741	1044357	58.83%	5.908%
49	17.127	2381	2387	2392	rVV7	6475	10770	0.61%	0.061%
50	17.215	2397	2402	2405	rVV5	1912	3003	0.17%	0.017%
51	17.309	2415	2418	2422	rBV4	1499	2418	0.14%	0.014%
52	17.392	2430	2432	2435	rBV3	1378	2224	0.13%	0.013%
53	17.456	2439	2443	2447	rVV3	2314	3907	0.22%	0.022%
54	17.598	2462	2467	2473	rVB2	41375	51559	2.90%	0.292%
55	17.662	2475	2478	2480	rBV3	2281	2973	0.17%	0.017%
56	17.698	2480	2484	2489	rVB6	2369	3842	0.22%	0.022%
57	17.821	2500	2505	2513	rBV	71534	100653	5.67%	0.569%
58	17.892	2515	2517	2520	rVB2	5377	5847	0.33%	0.033%
59	18.315	2587	2589	2592	rVB3	1873	2098	0.12%	0.012%
60	18.739	2656	2661	2662	rBV5	2742	4676	0.26%	0.026%
61	18.780	2664	2668	2675	rVB4	25983	32644	1.84%	0.185%
62	18.874	2680	2684	2689	rBV6	5277	7895	0.44%	0.045%
63	18.915	2689	2691	2692	rBV2	2747	2680	0.15%	0.015%
64	18.950	2693	2697	2700	rVV2	5665	7944	0.45%	0.045%
65	19.068	2715	2717	2720	rBV4	1977	2123	0.12%	0.012%
66	19.115	2721	2725	2730	rVV4	34385	47859	2.70%	0.271%
67	19.198	2736	2739	2740	rBV3	3839	3513	0.20%	0.020%
68	19.227	2740	2744	2749	rVB3	18626	22156	1.25%	0.125%
69	19.309	2752	2758	2767	rBV	1004472	1361519	76.69%	7.703%
70	20.844	3015	3019	3027	rBV2	89035	164407	9.26%	0.930%
71	21.109	3059	3064	3074	rBV	1354291	1775269	100.00%	10.044%
72	23.115	3399	3405	3411	rBV2	758699	1326552	74.72%	7.505%
73	23.250	3422	3428	3437	rVB2	936605	1719790	96.87%	9.730%

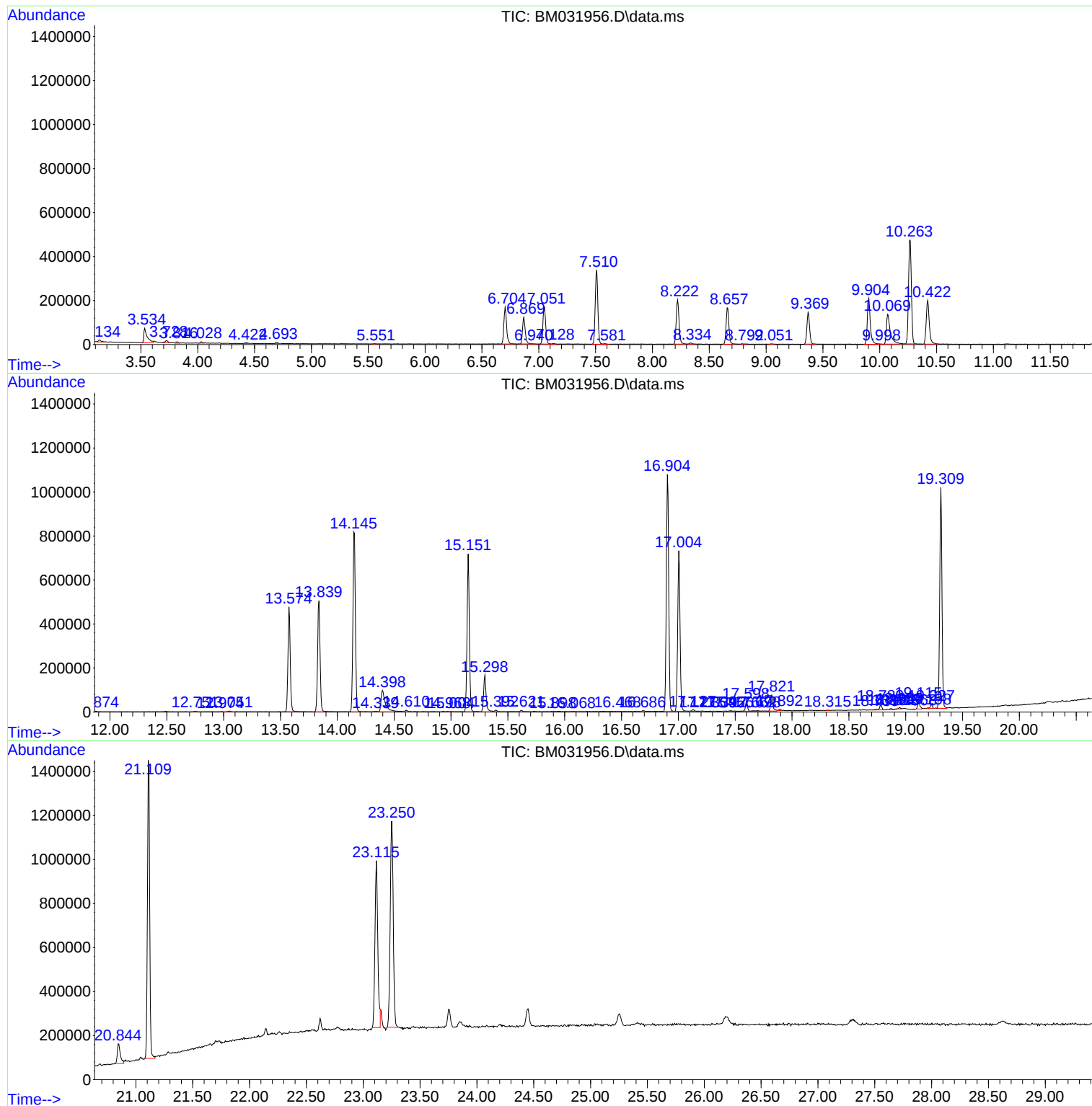
Sum of corrected areas: 17675731

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM082821.M
Quant Title : SVOA CALIBRATION

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



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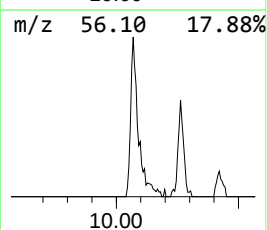
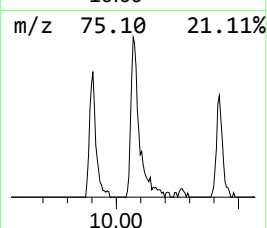
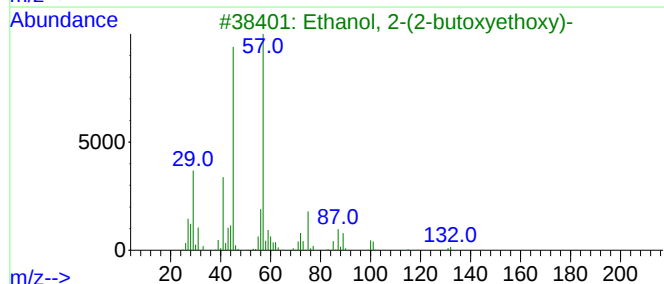
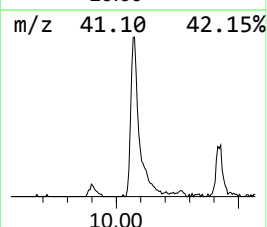
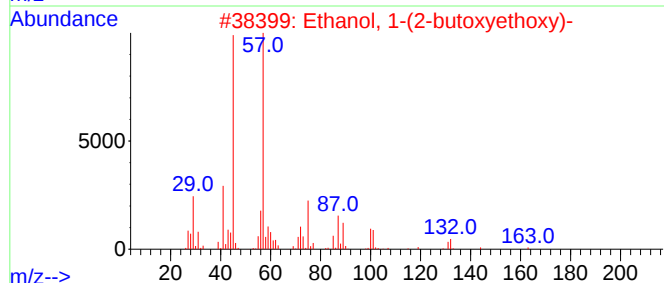
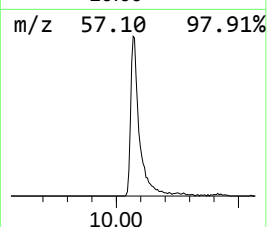
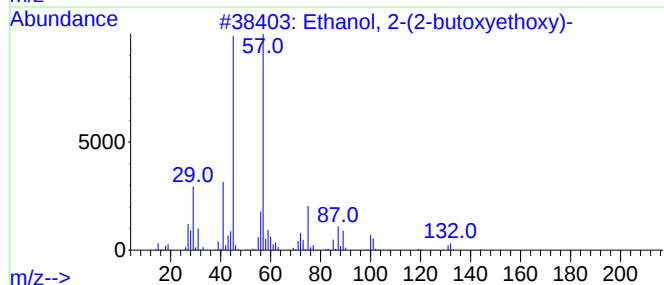
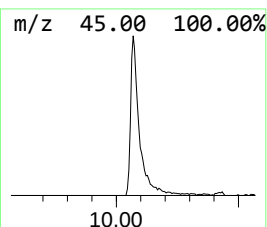
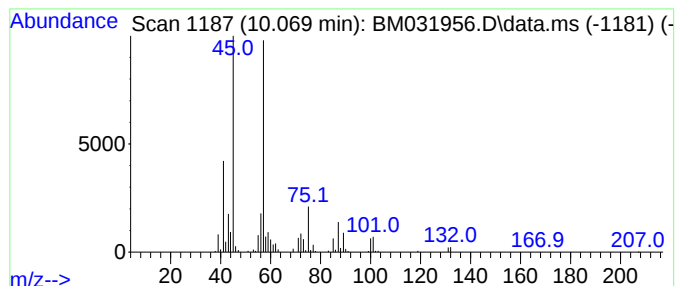
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM082821.M
Quant Title : SVOA CALIBRATION

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.
10.069	7.98 ng/ul	318972	Naphthalene-d8	10.269

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
2			Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	90
3			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
4			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
5			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	86



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TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Ethanol, 2-(2-b...	10.069	8.0	ng/u1	318972	2	10.269	799381	20.0