

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041823\
 Data File : BM039459.D
 Acq On : 18 Apr 2023 16:03
 Operator : CG/JU
 Sample : 02299-01
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AE3

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-BM041823.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BM039459.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.216	3	5	12	rVB	54642	65625	0.94%	0.111%
2	3.969	130	133	138	rVB	18921	25278	0.36%	0.043%
3	7.028	647	653	670	rBV	133030	304221	4.35%	0.517%
4	7.169	670	677	702	rVV	738939	1294763	18.53%	2.199%
5	7.369	705	711	735	rVV	675291	1217560	17.42%	2.068%
6	7.834	782	790	800	rBV	650950	1139636	16.31%	1.936%
7	8.569	908	915	929	rBV	475263	957454	13.70%	1.626%
8	8.945	974	979	982	rBV3	5872	8485	0.12%	0.014%
9	9.004	982	989	1009	rVV	875892	1597830	22.86%	2.714%
10	9.157	1010	1015	1035	rVB3	34478	108490	1.55%	0.184%
11	9.393	1052	1055	1060	rVB	7423	12828	0.18%	0.022%
12	9.728	1104	1112	1133	rBV	685330	1279328	18.31%	2.173%
13	10.275	1198	1205	1226	rBV	766837	1726468	24.70%	2.933%
14	10.639	1260	1267	1284	rBV	1047765	1832661	26.22%	3.113%
15	10.828	1297	1299	1305	rVB7	5505	7987	0.11%	0.014%
16	11.181	1354	1359	1365	rBV	124320	220419	3.15%	0.374%
17	11.245	1365	1370	1390	rVB	138993	332337	4.76%	0.565%
18	11.528	1413	1418	1419	rBV3	5731	8054	0.12%	0.014%
19	11.898	1475	1481	1487	rBV	12959	20902	0.30%	0.036%
20	12.245	1534	1540	1550	rBV2	15425	33689	0.48%	0.057%
21	12.845	1636	1642	1650	rBV4	7669	16977	0.24%	0.029%
22	13.098	1680	1685	1691	rBV2	12899	19117	0.27%	0.032%
23	13.904	1816	1822	1837	rBV	2543642	3721297	53.25%	6.321%
24	14.186	1862	1870	1885	rBV	2684942	4195226	60.03%	7.126%
25	14.492	1914	1922	1941	rBV	1767771	2687494	38.45%	4.565%
26	14.763	1961	1968	1969	rBV2	24665	39419	0.56%	0.067%
27	15.280	2054	2056	2059	rBV	6224	9209	0.13%	0.016%
28	15.316	2059	2062	2068	rVB	14440	21789	0.31%	0.037%
29	15.486	2082	2091	2108	rBV	3600711	5389863	77.12%	9.156%
30	15.622	2108	2114	2129	rVV	894060	1492301	21.35%	2.535%
31	15.727	2129	2132	2139	rVB4	21245	37861	0.54%	0.064%
32	15.857	2149	2154	2160	rBV4	12281	18426	0.26%	0.031%
33	16.704	2292	2298	2300	rBV7	4333	7325	0.10%	0.012%
34	16.839	2317	2321	2325	rVB5	6545	10125	0.14%	0.017%
35	17.039	2352	2355	2359	rVB4	7874	12322	0.18%	0.021%
36	17.163	2372	2376	2379	rBV5	5458	8688	0.12%	0.015%

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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-BM041823.MA.M
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37	17.245	2383	2390	2400	rVV	2177966	3211220	45.95%	5.455%
38	17.339	2400	2406	2433	rVV	3871922	6043043	86.47%	10.265%
39	17.527	2435	2438	2447	rVB	23337	36008	0.52%	0.061%
40	17.910	2498	2503	2508	rBV	170955	221546	3.17%	0.376%
41	17.951	2508	2510	2518	rVV9	9062	15833	0.23%	0.027%
42	18.027	2521	2523	2529	rVB6	5472	7525	0.11%	0.013%
43	18.145	2539	2543	2547	rBV2	45301	84303	1.21%	0.143%
44	18.215	2551	2555	2566	rVB	54679	98646	1.41%	0.168%
45	19.204	2719	2723	2730	rBV	52240	82569	1.18%	0.140%
46	19.280	2732	2736	2743	rVV	41041	71010	1.02%	0.121%
47	19.427	2757	2761	2774	rBV7	37576	109237	1.56%	0.186%
48	19.568	2782	2785	2791	rBV5	16853	21052	0.30%	0.036%
49	19.639	2791	2797	2808	rBV	4982026	6988726	100.00%	11.872%
50	21.439	3098	3103	3113	rBV	2287105	3213226	45.98%	5.458%
51	23.656	3473	3480	3496	rVB2	2752241	5924412	84.77%	10.064%
52	23.809	3500	3506	3520	rVB2	1390572	2859004	40.91%	4.857%

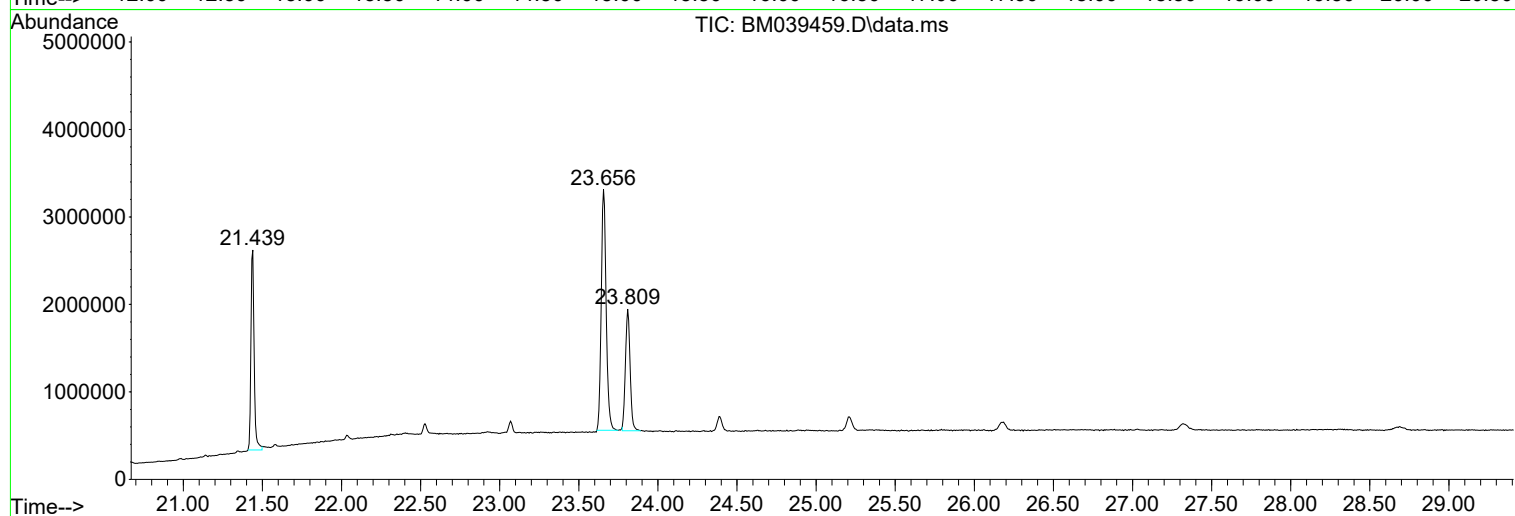
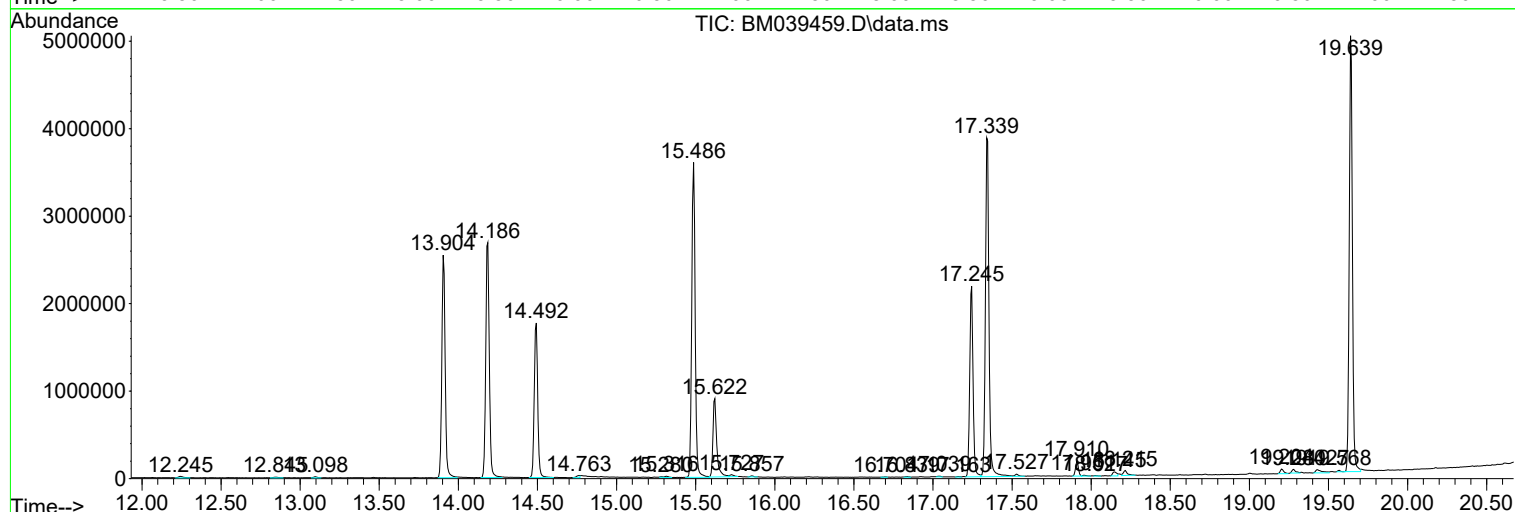
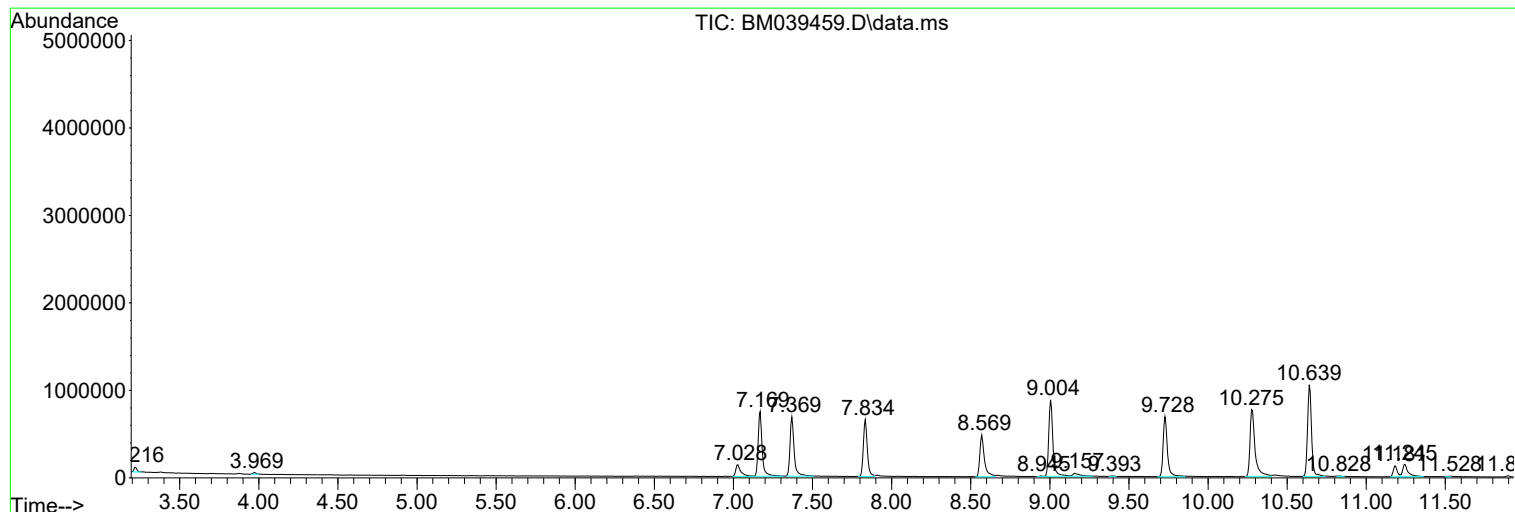
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM041823.MA.M
Quant Title : SVOA CALIBRATION

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



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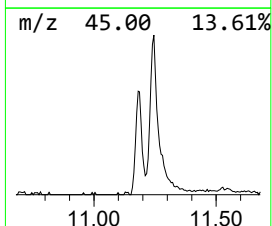
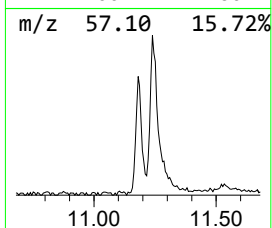
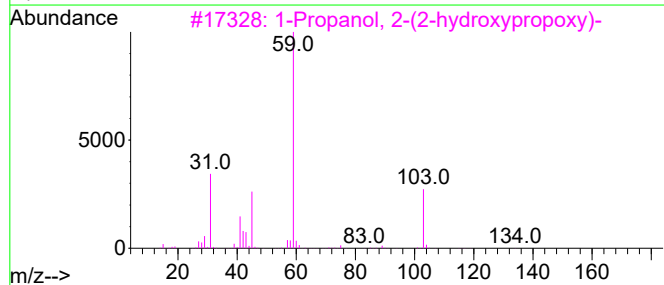
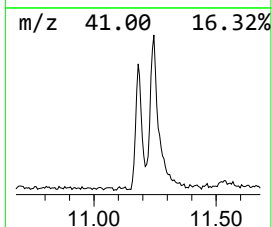
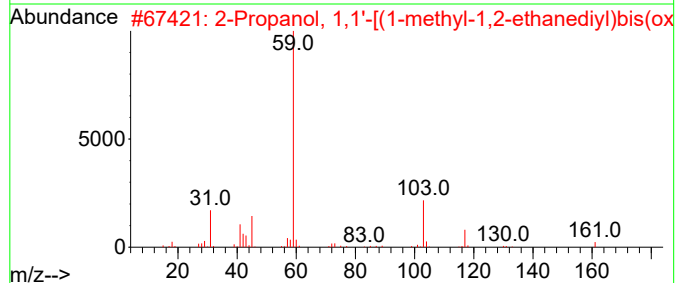
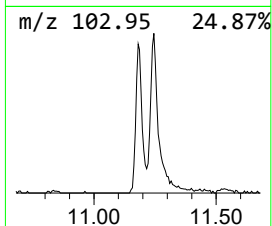
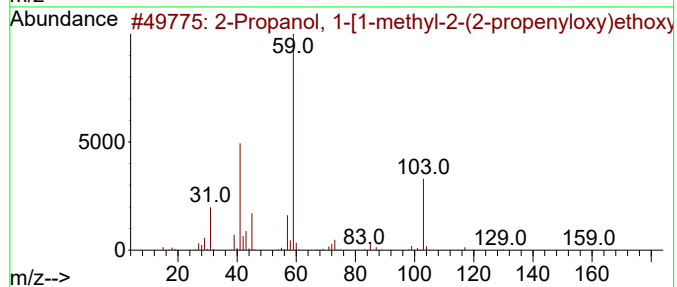
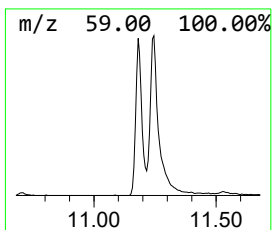
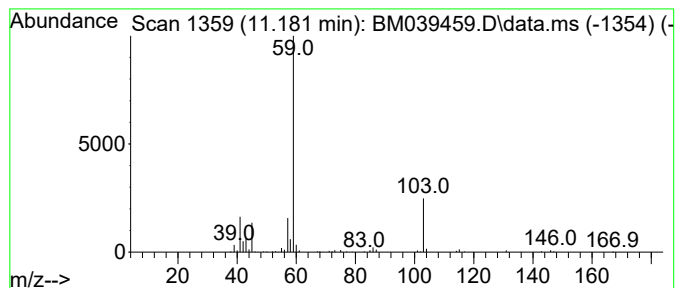
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM041823.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 2-Propanol, 1-[1-methyl-2-(... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.181	2.41 ng/ul	220419	Naphthalene-d8	10.640

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanol, 1-[1-methyl-2-(2-pro...	174	C9H18O3	055956-25-7	83		
2	2-Propanol, 1,1'-[(1-methyl-1,2-...	192	C9H20O4	001638-16-0	78		
3	1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	64		
4	1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	64		
5	1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	64		



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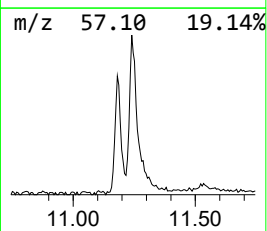
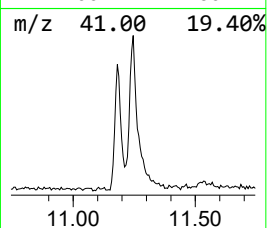
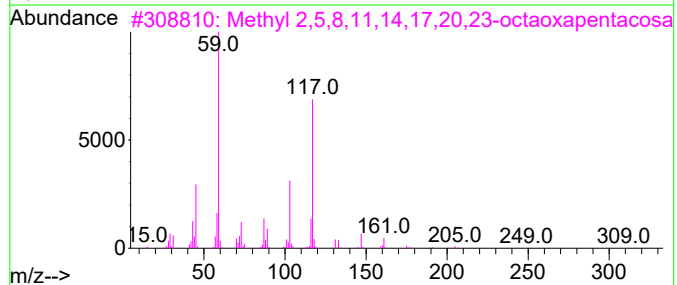
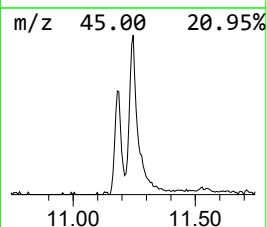
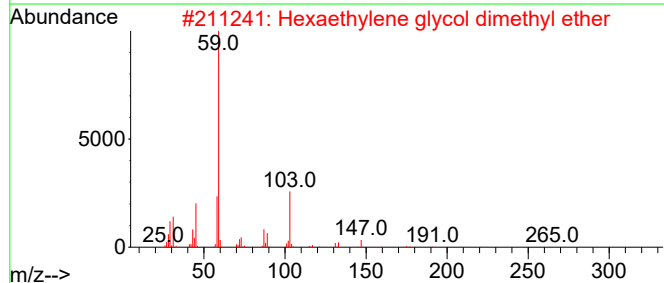
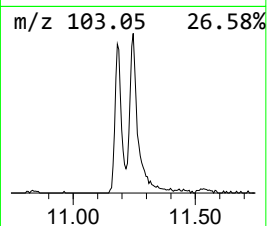
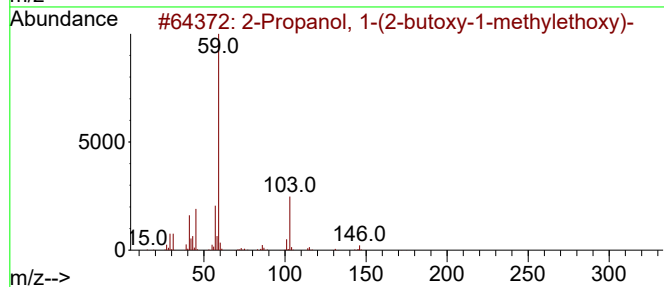
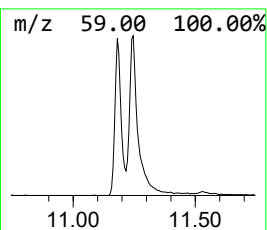
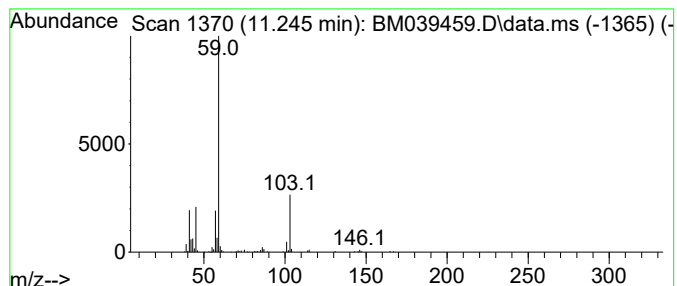
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM041823.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 2-Propanol, 1-(2-butoxy-1-m... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.245	3.63 ng/ul	332337	Naphthalene-d8	10.640

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Propanol, 1-(2-butoxy-1-methyl...	190	C10H22O3	029911-28-2	90
2			Hexaethylene glycol dimethyl ether	310	C14H30O7	001072-40-8	78
3			Methyl 2,5,8,11,14,17,20,23-octa...	412	C18H36O10	1000366-81-2	74
4			2-Propanol, 1-[1-methyl-2-(2-pro...	174	C9H18O3	055956-25-7	64
5			Methyl 2,5,8,11,14,17,20-heptaox...	368	C16H32O9	1010366-81-3	64



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Quant Title : SVOA CALIBRATION

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-Propanol, 1-[... 11.181	2.4	ng/ul	220419	2	10.640	1832660	20.0		
2-Propanol, 1-(... 11.245	3.6	ng/ul	332337	2	10.640	1832660	20.0		