

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092322\
 Data File : BM036683.D
 Acq On : 23 Sep 2022 18:47
 Operator : CG/JU
 Sample : N4706-10
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 CB8L9

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-BM091522.M
 Title : SVOA CALIBRATION

Signal : TIC: BM036683.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.123	3	6	34	rVB	9068060	12685187	92.01%	8.959%
2	3.393	48	52	57	rBV	105954	139600	1.01%	0.099%
3	3.675	97	100	110	rBV	58873	89816	0.65%	0.063%
4	3.823	122	125	139	rBV	643369	970294	7.04%	0.685%
5	4.058	160	165	175	rVB	64706	114257	0.83%	0.081%
6	4.152	177	181	187	rVB	33286	48411	0.35%	0.034%
7	4.605	255	258	267	rVB2	39776	71135	0.52%	0.050%
8	5.093	337	341	345	rBV2	13173	16688	0.12%	0.012%
9	5.305	372	377	381	rBV	16716	29144	0.21%	0.021%
10	5.699	440	444	449	rVB4	10036	16233	0.12%	0.011%
11	5.752	449	453	459	rBV4	13949	24599	0.18%	0.017%
12	6.599	593	597	602	rVB7	9595	17869	0.13%	0.013%
13	7.005	660	666	671	rBV9	7599	14734	0.11%	0.010%
14	7.175	687	695	707	rVB	628750	1011589	7.34%	0.714%
15	7.346	717	724	733	rBV	1878088	2928127	21.24%	2.068%
16	7.546	749	758	767	rBV	1923746	3044523	22.08%	2.150%
17	7.640	770	774	781	rVB9	12170	26513	0.19%	0.019%
18	8.016	831	838	845	rBV	1800780	2811157	20.39%	1.985%
19	8.087	845	850	854	rVB5	10270	17394	0.13%	0.012%
20	8.728	949	959	969	rBV	1734594	2854338	20.70%	2.016%
21	9.010	1002	1007	1012	rBV	25913	42126	0.31%	0.030%
22	9.181	1028	1036	1047	rBV	2262648	3601665	26.12%	2.544%
23	9.299	1051	1056	1060	rVV6	11319	19064	0.14%	0.013%
24	9.352	1060	1065	1070	rVB8	9292	14685	0.11%	0.010%
25	9.546	1091	1098	1103	rBV10	21848	47055	0.34%	0.033%
26	9.604	1105	1108	1112	rVB4	10243	17043	0.12%	0.012%
27	9.904	1149	1159	1169	rBV	1732023	2789373	20.23%	1.970%
28	10.134	1192	1198	1202	rVB8	8682	15874	0.12%	0.011%
29	10.328	1226	1231	1236	rBV7	11608	23038	0.17%	0.016%
30	10.446	1243	1251	1261	rBV	2848288	4657144	33.78%	3.289%
31	10.604	1274	1278	1287	rVB	51633	93454	0.68%	0.066%
32	10.828	1308	1316	1325	rVB	2659096	4361142	31.63%	3.080%
33	10.963	1330	1339	1355	rBV	2147247	3545935	25.72%	2.504%
34	11.193	1374	1378	1382	rBV6	9190	16374	0.12%	0.012%
35	11.487	1423	1428	1432	rBV6	8968	14123	0.10%	0.010%
36	11.616	1444	1450	1454	rBV5	18783	33220	0.24%	0.023%

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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 >

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
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37	11.734	1464	1470	1477	rBV10	10248	20775	0.15%	0.015%
38	12.163	1539	1543	1547	rVB	19154	30741	0.22%	0.022%
39	12.322	1563	1570	1576	rBV5	19865	36287	0.26%	0.026%
40	12.404	1578	1584	1590	rBV2	54643	111874	0.81%	0.079%
41	13.016	1679	1688	1693	rBV8	16602	35373	0.26%	0.025%
42	13.263	1726	1730	1735	rVB7	15636	24309	0.18%	0.017%
43	13.375	1746	1749	1754	rBV5	7429	14214	0.10%	0.010%
44	13.545	1774	1778	1782	rBV6	11679	15439	0.11%	0.011%
45	14.004	1846	1856	1857	rBV6	16810	35619	0.26%	0.025%
46	14.057	1858	1865	1880	rVB	5479577	7452683	54.06%	5.263%
47	14.287	1900	1904	1906	rBV5	16482	24352	0.18%	0.017%
48	14.334	1906	1912	1927	rVB	5999192	8826814	64.03%	6.234%
49	14.569	1948	1952	1957	rVB4	29056	36894	0.27%	0.026%
50	14.639	1957	1964	1971	rBV	4278790	5973871	43.33%	4.219%
51	14.704	1971	1975	1981	rVB	59640	93710	0.68%	0.066%
52	14.834	1990	1997	2007	rBV	599787	913789	6.63%	0.645%
53	14.992	2021	2024	2030	rBV7	12147	21335	0.15%	0.015%
54	15.051	2031	2034	2038	rVB5	23817	31834	0.23%	0.022%
55	15.381	2085	2090	2094	rBV	144629	193719	1.41%	0.137%
56	15.434	2096	2099	2106	rVB2	21788	39248	0.28%	0.028%
57	15.628	2125	2132	2139	rBV	8196951	11316861	82.09%	7.992%
58	15.739	2145	2151	2160	rBV	2164361	2915099	21.14%	2.059%
59	15.863	2168	2172	2180	rVB3	73385	140101	1.02%	0.099%
60	16.139	2214	2219	2224	rBV	166443	229232	1.66%	0.162%
61	16.410	2261	2265	2269	rBV5	24938	35718	0.26%	0.025%
62	16.463	2271	2274	2278	rBV2	19633	28198	0.20%	0.020%
63	16.645	2301	2305	2310	rBV	83673	110636	0.80%	0.078%
64	17.375	2423	2429	2438	rBV	4943149	6929881	50.27%	4.894%
65	17.480	2440	2447	2457	rVV	8310868	12216560	88.61%	8.628%
66	18.045	2541	2543	2548	rVB2	46634	54062	0.39%	0.038%
67	18.269	2577	2581	2588	rBV2	263587	408959	2.97%	0.289%
68	19.327	2757	2761	2766	rBV	103047	145278	1.05%	0.103%
69	19.392	2769	2772	2775	rBV	110119	137390	1.00%	0.097%
70	19.539	2794	2797	2802	rBV2	140223	191730	1.39%	0.135%
71	19.763	2829	2835	2840	rBV	10319163	13786290	100.00%	9.736%
72	20.757	3001	3004	3007	rBV	220645	216579	1.57%	0.153%
73	21.251	3085	3088	3096	rBV2	348846	445955	3.23%	0.315%
74	21.545	3133	3138	3145	rBV	5182694	6511062	47.23%	4.598%
75	23.833	3520	3527	3534	rBV2	5192594	9948650	72.16%	7.026%
76	23.986	3547	3553	3563	rVB2	2844240	5669875	41.13%	4.004%

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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_M\Methods\SFAM-EPA-BM091522.M
Title : SVOA CALIBRATION

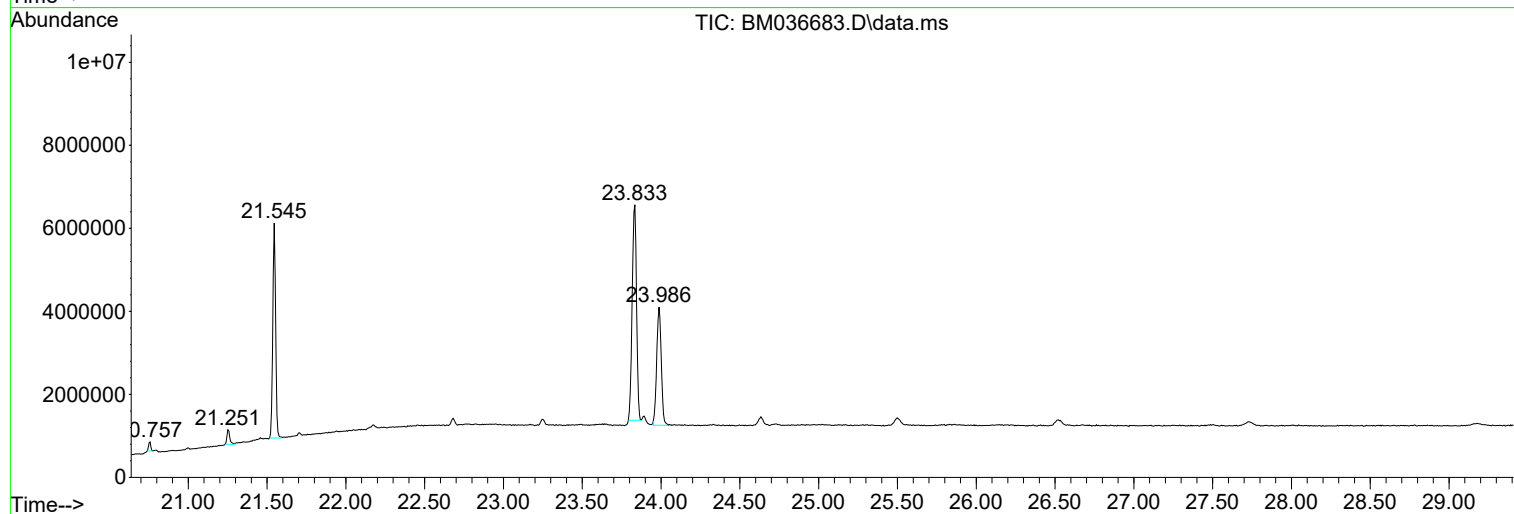
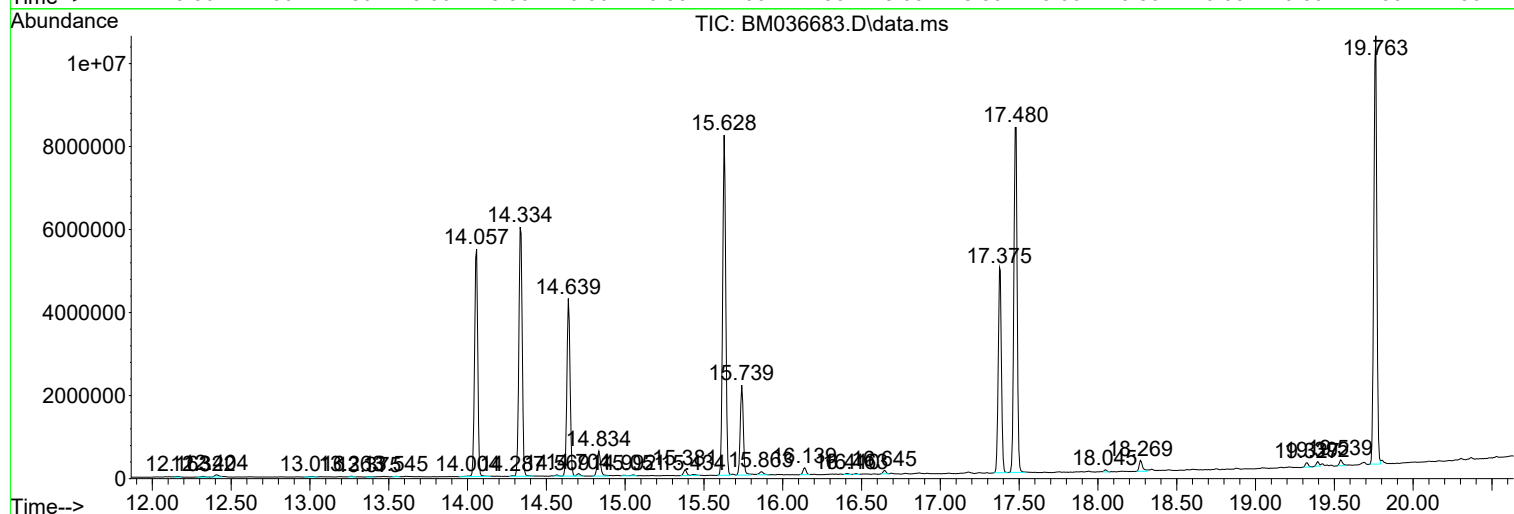
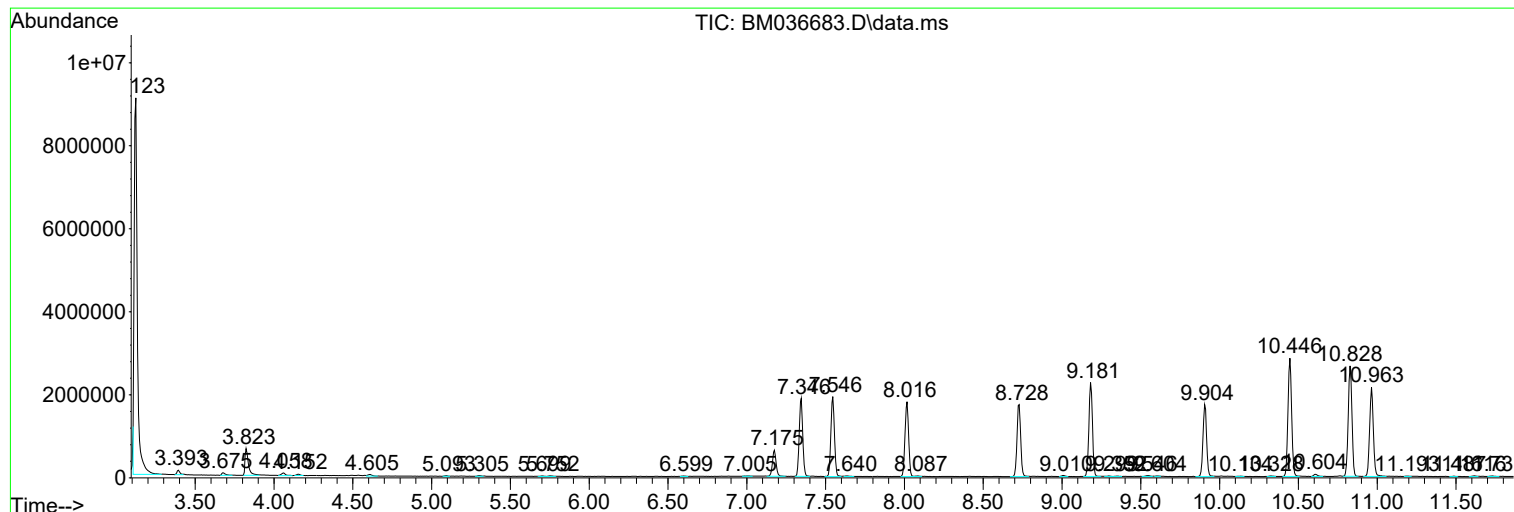
Sum of corrected areas: 141593919

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Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
Quant Title : SVOA CALIBRATION

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



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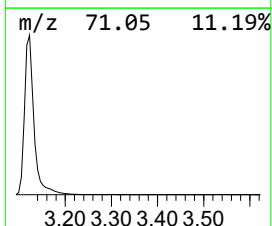
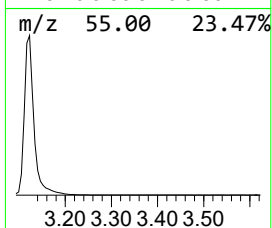
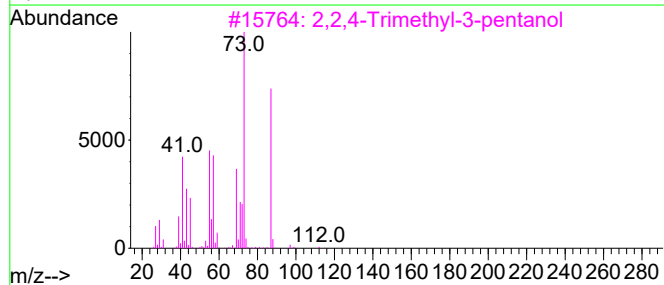
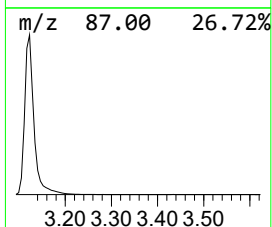
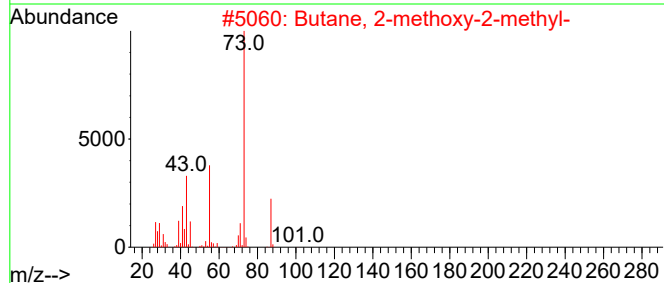
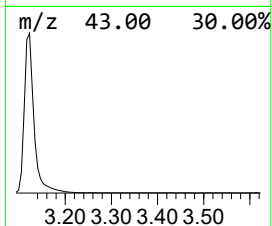
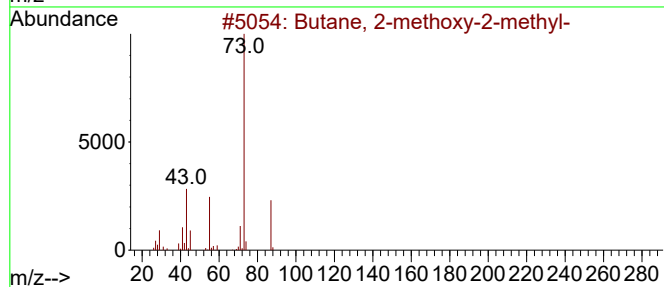
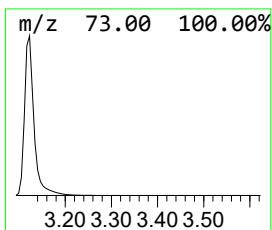
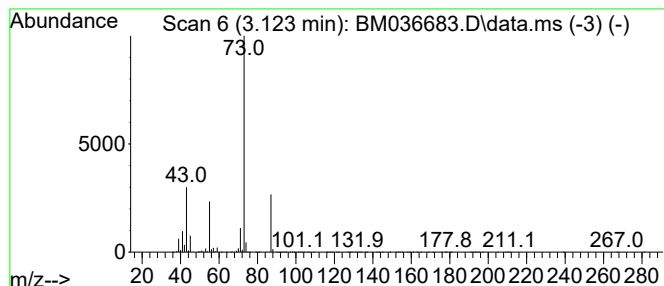
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.123	90.25 ng/ul	12685200	1,4-Dichlorobenzene-d4	8.016

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
3			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	43
4			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
5			Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9



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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
Butane, 2-metho...	3.123	90.3	ng/ul	12685200	1	8.016	2811160	20.0