

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092322\
 Data File : BM036728.D
 Acq On : 25 Sep 2022 00:12
 Operator : CG/JU
 Sample : N4662-06
 Misc :
 ALS Vial : 63 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 EXZ27

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: BM036729.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.110	3	4	26	rVB	5836699	7074893	77.87%	7.909%
2	3.387	47	51	54	rBV	64648	90516	1.00%	0.101%
3	3.422	54	57	64	rVB	34479	54099	0.60%	0.060%
4	3.822	122	125	136	rBV	155299	271525	2.99%	0.304%
5	4.051	159	164	172	rVB	41673	63730	0.70%	0.071%
6	4.146	175	180	188	rVB	71094	104088	1.15%	0.116%
7	4.704	269	275	280	rVB9	5381	12167	0.13%	0.014%
8	5.416	392	396	399	rBV6	7583	12346	0.14%	0.014%
9	5.487	405	408	412	rVB2	9464	14579	0.16%	0.016%
10	5.569	417	422	425	rVV4	7494	11222	0.12%	0.013%
11	5.622	427	431	439	rBV2	16724	35230	0.39%	0.039%
12	5.793	456	460	464	rVB6	5808	10004	0.11%	0.011%
13	5.998	490	495	501	rVB3	38141	60531	0.67%	0.068%
14	6.263	536	540	545	rBV8	8249	14981	0.16%	0.017%
15	6.481	570	577	578	rBV7	8118	10066	0.11%	0.011%
16	6.551	584	589	594	rBV5	8850	15082	0.17%	0.017%
17	6.622	598	601	603	rBV4	9062	11343	0.12%	0.013%
18	6.981	658	662	668	rVV7	12210	22979	0.25%	0.026%
19	7.039	670	672	677	rVV5	9649	17358	0.19%	0.019%
20	7.092	677	681	686	rVV5	16555	34177	0.38%	0.038%
21	7.157	686	692	707	rVB	334338	595854	6.56%	0.666%
22	7.334	715	722	729	rBV	1296959	2047723	22.54%	2.289%
23	7.534	749	756	766	rVV	1200158	1894613	20.85%	2.118%
24	7.622	766	771	783	rVB2	32828	70924	0.78%	0.079%
25	8.004	827	836	843	rBV	1284626	2060855	22.68%	2.304%
26	8.069	843	847	850	rVB5	5497	9219	0.10%	0.010%
27	8.292	879	885	890	rBV8	6730	9891	0.11%	0.011%
28	8.557	924	930	941	rVB	82206	143874	1.58%	0.161%
29	8.710	950	956	964	rBV	1034619	1615943	17.79%	1.806%
30	9.169	1027	1034	1043	rVV	1553174	2495696	27.47%	2.790%
31	9.233	1043	1045	1050	rVB4	14845	19956	0.22%	0.022%
32	9.328	1057	1061	1065	rBV7	5482	11206	0.12%	0.013%
33	9.892	1147	1157	1170	rBV	1098695	1802647	19.84%	2.015%
34	9.980	1170	1172	1177	rVB6	5596	9595	0.11%	0.011%
35	10.433	1240	1249	1259	rBV	1723622	2876503	31.66%	3.216%
36	10.598	1272	1277	1282	rBV3	24261	45174	0.50%	0.050%

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Min Area: 0.1 % of largest Peak

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Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

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Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
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37	10.816	1303	1314	1325	rBV	1884629	3157320	34.75%	3.530%
38	10.951	1330	1337	1349	rBV	919037	1562409	17.20%	1.747%
39	11.598	1440	1447	1450	rBV9	4928	9349	0.10%	0.010%
40	12.186	1544	1547	1553	rBV8	7915	13681	0.15%	0.015%
41	12.392	1576	1582	1590	rVB	37364	60280	0.66%	0.067%
42	12.669	1625	1629	1636	rVB10	13937	23668	0.26%	0.026%
43	13.257	1723	1729	1737	rVB	15755	29218	0.32%	0.033%
44	13.457	1759	1763	1767	rVV6	7510	12918	0.14%	0.014%
45	13.533	1770	1776	1784	rVV	51982	86286	0.95%	0.096%
46	13.604	1784	1788	1793	rVB7	6997	11999	0.13%	0.013%
47	13.804	1816	1822	1827	rVB9	5983	11557	0.13%	0.013%
48	14.039	1854	1862	1872	rBV	3451166	4713378	51.88%	5.269%
49	14.321	1903	1910	1923	rBV	3784927	5614914	61.80%	6.277%
50	14.521	1941	1944	1948	rVB6	7205	10732	0.12%	0.012%
51	14.627	1955	1962	1973	rBV	2873535	4184041	46.05%	4.677%
52	14.815	1989	1994	2008	rBV	230559	390070	4.29%	0.436%
53	14.986	2018	2023	2028	rBV9	5588	10705	0.12%	0.012%
54	15.045	2028	2033	2040	rVB9	16897	27096	0.30%	0.030%
55	15.615	2123	2130	2142	rBV	5381591	7564585	83.26%	8.456%
56	15.727	2142	2149	2161	rBV	1216627	1645738	18.11%	1.840%
57	15.857	2166	2171	2176	rVB	46506	72673	0.80%	0.081%
58	16.174	2223	2225	2231	rVB7	5787	9506	0.10%	0.011%
59	16.304	2239	2247	2250	rBV9	8455	18978	0.21%	0.021%
60	16.398	2259	2263	2267	rVB5	13319	19002	0.21%	0.021%
61	16.851	2335	2340	2345	rBV4	12484	25996	0.29%	0.029%
62	17.045	2369	2373	2377	rBV6	9340	13214	0.15%	0.015%
63	17.368	2421	2428	2438	rBV	3311237	4723785	51.99%	5.281%
64	17.462	2438	2444	2456	rBV	6281735	8622307	94.90%	9.639%
65	17.798	2498	2501	2507	rVB	21083	30347	0.33%	0.034%
66	17.921	2519	2522	2529	rVB3	12024	16085	0.18%	0.018%
67	18.033	2539	2541	2545	rVB5	9453	10219	0.11%	0.011%
68	18.256	2575	2579	2587	rBV	74430	107104	1.18%	0.120%
69	19.315	2755	2759	2763	rBV2	58471	78215	0.86%	0.087%
70	19.386	2767	2771	2774	rBV	67943	89445	0.98%	0.100%
71	19.527	2792	2795	2800	rBV5	48214	65977	0.73%	0.074%
72	19.745	2827	2832	2840	rBV	6841845	9085881	100.00%	10.157%
73	20.739	2999	3001	3005	rVB2	119580	112539	1.24%	0.126%
74	21.533	3131	3136	3142	rBV	3032553	3910105	43.03%	4.371%
75	23.809	3516	3523	3530	rBV2	3285602	6295007	69.28%	7.037%
76	23.962	3543	3549	3558	rVB2	1643646	3357342	36.95%	3.753%

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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

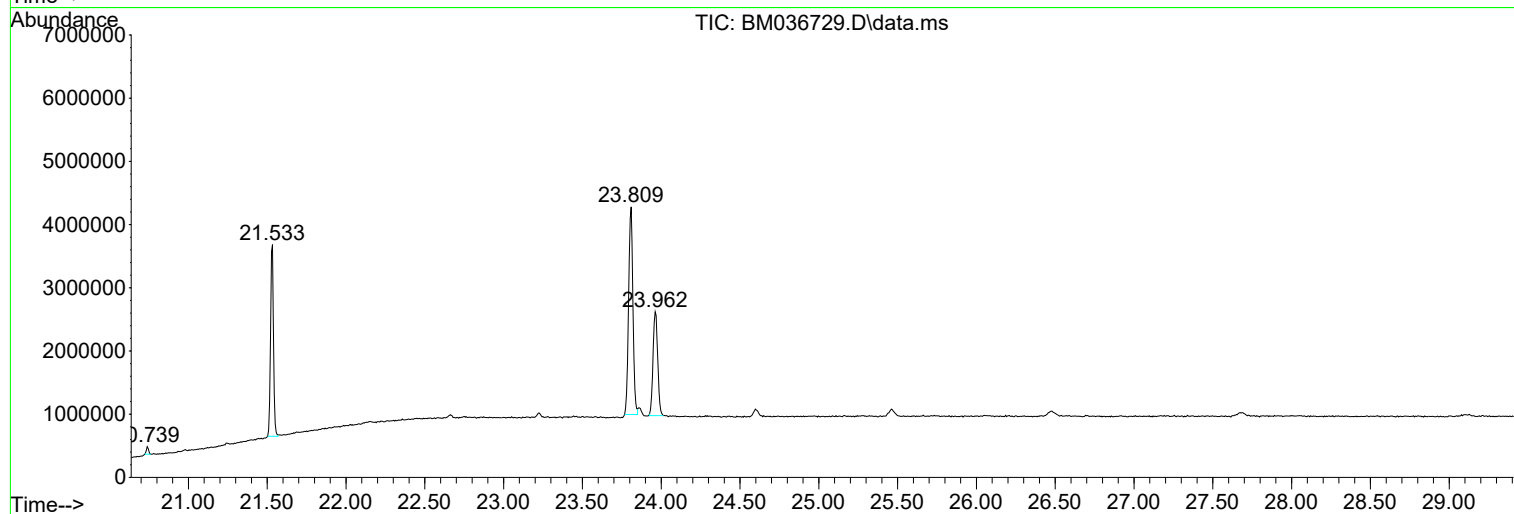
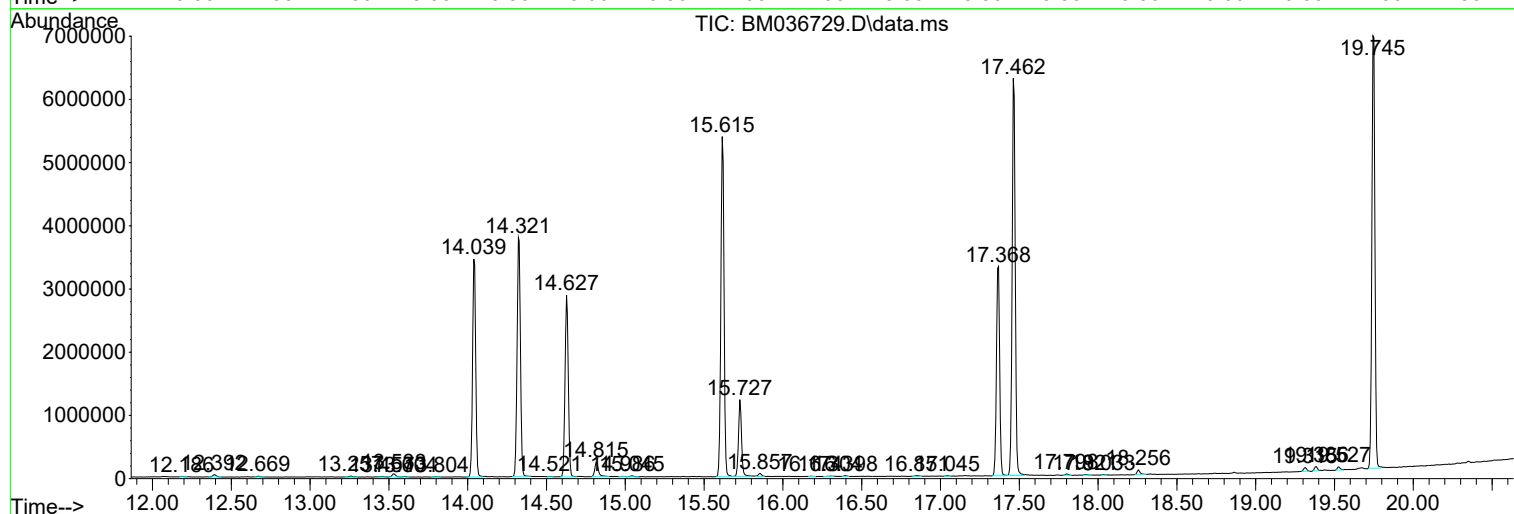
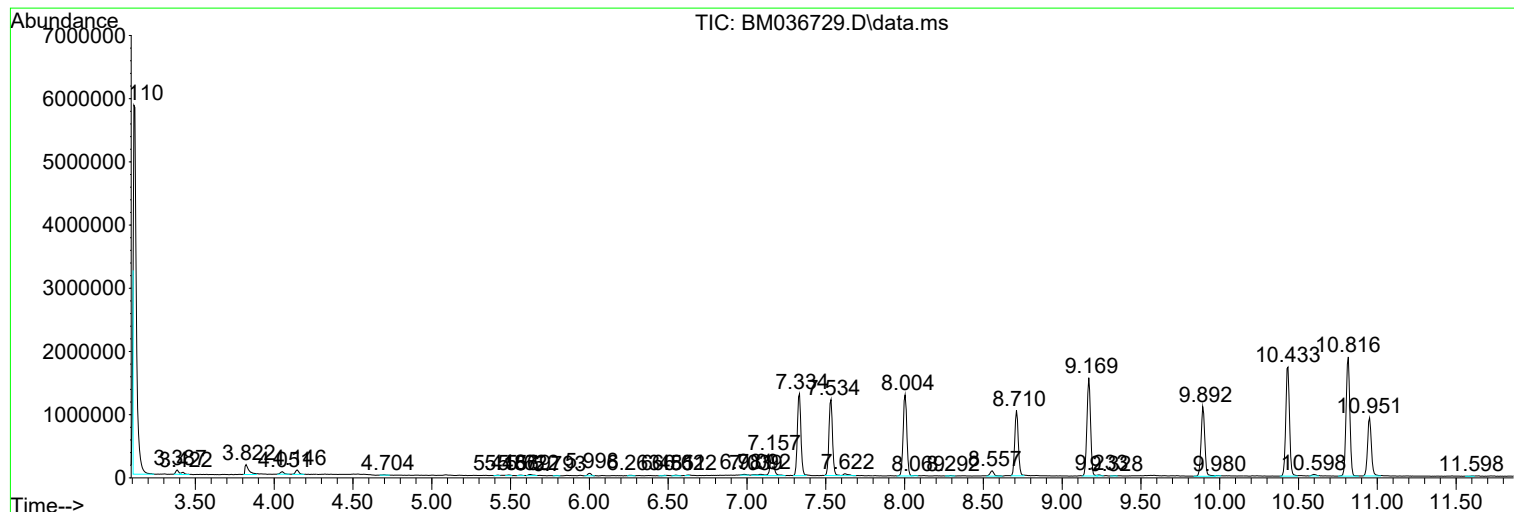
Sum of corrected areas: 89454260

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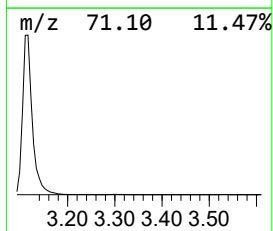
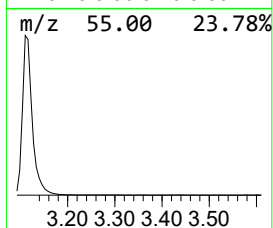
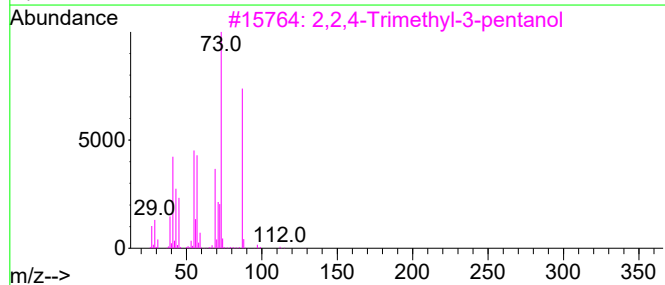
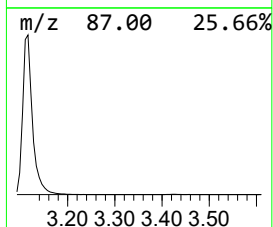
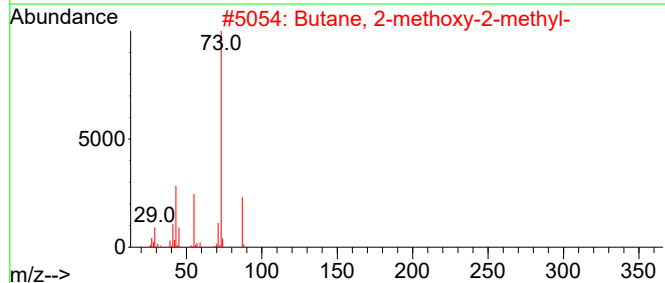
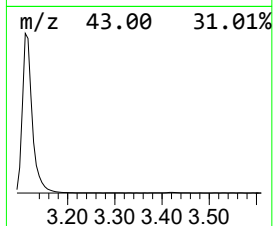
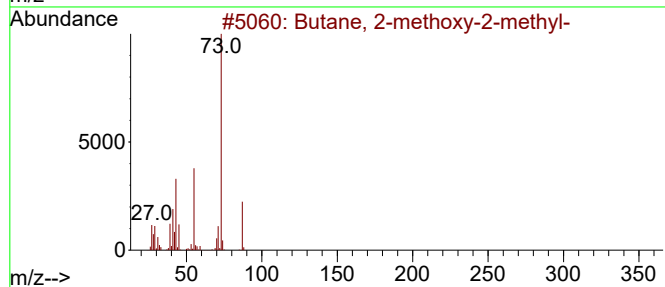
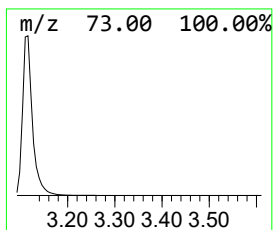
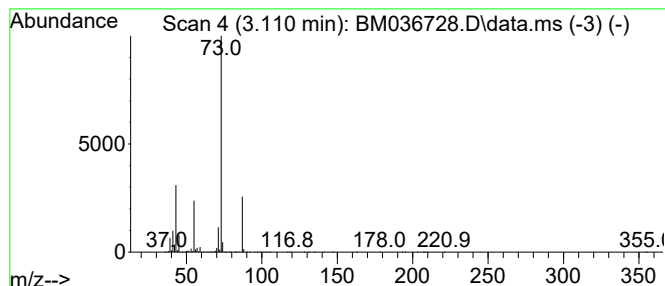
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.110	68.66 ng/ul	7074890	1,4-Dichlorobenzene-d4	8.004

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	39
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...	3.110	68.7	ng/ul	7074890	1	8.004	2060860	20.0