

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042522\
 Data File : BP010082.D
 Acq On : 25 Apr 2022 14:51
 Operator : CG/JU
 Sample : N2475-03
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A0BM3

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_P\Methods\SFAM-EPA-BP042122.M
 Title : SVOA CALIBRATION

Signal : TIC: BP010082.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.111	3	4	17	rVB	1049867	915898	17.92%	2.242%
2	3.369	44	48	54	rBV	48779	60561	1.18%	0.148%
3	3.793	117	120	132	rBV	202142	330759	6.47%	0.810%
4	4.122	172	176	181	rBV	28297	40217	0.79%	0.098%
5	4.452	228	232	237	rBV3	5964	7058	0.14%	0.017%
6	5.046	327	333	350	rBV	3706377	5111993	100.00%	12.515%
7	6.234	529	535	537	rBV7	3859	7260	0.14%	0.018%
8	7.105	677	683	691	rBV	74672	124598	2.44%	0.305%
9	7.269	704	711	721	rBV	571826	932485	18.24%	2.283%
10	7.475	739	746	757	rBV	263718	446729	8.74%	1.094%
11	7.569	760	762	771	rVB8	4748	9006	0.18%	0.022%
12	7.940	816	825	833	rBV	591702	984571	19.26%	2.410%
13	8.010	834	837	844	rVB	12863	21205	0.41%	0.052%
14	8.646	939	945	958	rBV	247669	442859	8.66%	1.084%
15	8.763	962	965	974	rVB5	10777	22485	0.44%	0.055%
16	9.099	1014	1022	1035	rBV	687042	1194230	23.36%	2.924%
17	9.828	1137	1146	1156	rBV	152651	264292	5.17%	0.647%
18	10.369	1230	1238	1249	rBV	330879	594329	11.63%	1.455%
19	10.746	1295	1302	1317	rBV	751936	1336879	26.15%	3.273%
20	10.881	1317	1325	1346	rBV	677159	1291737	25.27%	3.162%
21	12.093	1528	1531	1537	rVB7	4662	7608	0.15%	0.019%
22	12.157	1537	1542	1546	rBV7	4717	8086	0.16%	0.020%
23	12.287	1560	1564	1567	rBV6	3706	5410	0.11%	0.013%
24	12.340	1567	1573	1579	rVV3	14458	24301	0.48%	0.059%
25	13.316	1734	1739	1744	rVB9	2355	5248	0.10%	0.013%
26	13.481	1760	1767	1774	rBV	59130	97950	1.92%	0.240%
27	13.634	1789	1793	1801	rVB8	6599	14140	0.28%	0.035%
28	13.898	1829	1838	1844	rBV5	22184	41300	0.81%	0.101%
29	13.981	1846	1852	1868	rBV	1549209	2268129	44.37%	5.553%
30	14.145	1878	1880	1887	rVB8	3027	5441	0.11%	0.013%
31	14.269	1890	1901	1918	rBV	1654023	2514341	49.19%	6.155%
32	14.410	1920	1925	1929	rBV7	5500	11261	0.22%	0.028%
33	14.575	1945	1953	1967	rBV	1215656	1905125	37.27%	4.664%
34	14.798	1985	1991	2002	rBV3	26282	54338	1.06%	0.133%
35	15.392	2086	2092	2098	rBV2	16136	25039	0.49%	0.061%
36	15.569	2115	2122	2136	rBV	2282304	3381851	66.16%	8.279%

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37	15.687	2136	2142	2158	rVV	156319	280158	5.48%	0.686%
38	15.810	2158	2163	2171	rVV	29133	51120	1.00%	0.125%
39	15.875	2171	2174	2178	rVV2	11256	16097	0.31%	0.039%
40	15.939	2181	2185	2190	rVB8	3847	6487	0.13%	0.016%
41	16.269	2235	2241	2243	rBV7	3139	5171	0.10%	0.013%
42	16.357	2251	2256	2261	rVB6	8829	14482	0.28%	0.035%
43	16.834	2334	2337	2344	rVB9	3821	5621	0.11%	0.014%
44	16.992	2358	2364	2365	rBV6	5600	8845	0.17%	0.022%
45	17.075	2374	2378	2383	rBV7	4959	7869	0.15%	0.019%
46	17.328	2414	2421	2431	rBV	1337446	1957613	38.29%	4.792%
47	17.428	2431	2438	2459	rVB2	2602238	3971684	77.69%	9.723%
48	17.616	2467	2470	2477	rVB5	5624	8743	0.17%	0.021%
49	18.110	2551	2554	2558	rVB6	5942	8226	0.16%	0.020%
50	18.216	2568	2572	2579	rBV8	6972	15809	0.31%	0.039%
51	18.280	2579	2583	2588	rVV	17321	21851	0.43%	0.053%
52	18.722	2654	2658	2666	rVB9	6693	10921	0.21%	0.027%
53	19.239	2740	2746	2753	rBV4	33649	58630	1.15%	0.144%
54	19.304	2753	2757	2764	rVV	28687	51229	1.00%	0.125%
55	19.657	2811	2817	2834	rBV	2938749	4203232	82.22%	10.290%
56	21.410	3110	3115	3129	rBV	1027004	1607599	31.45%	3.936%
57	23.710	3499	3506	3514	rBV	1262633	2733690	53.48%	6.692%
58	23.874	3527	3534	3546	rVB	572090	1293754	25.31%	3.167%

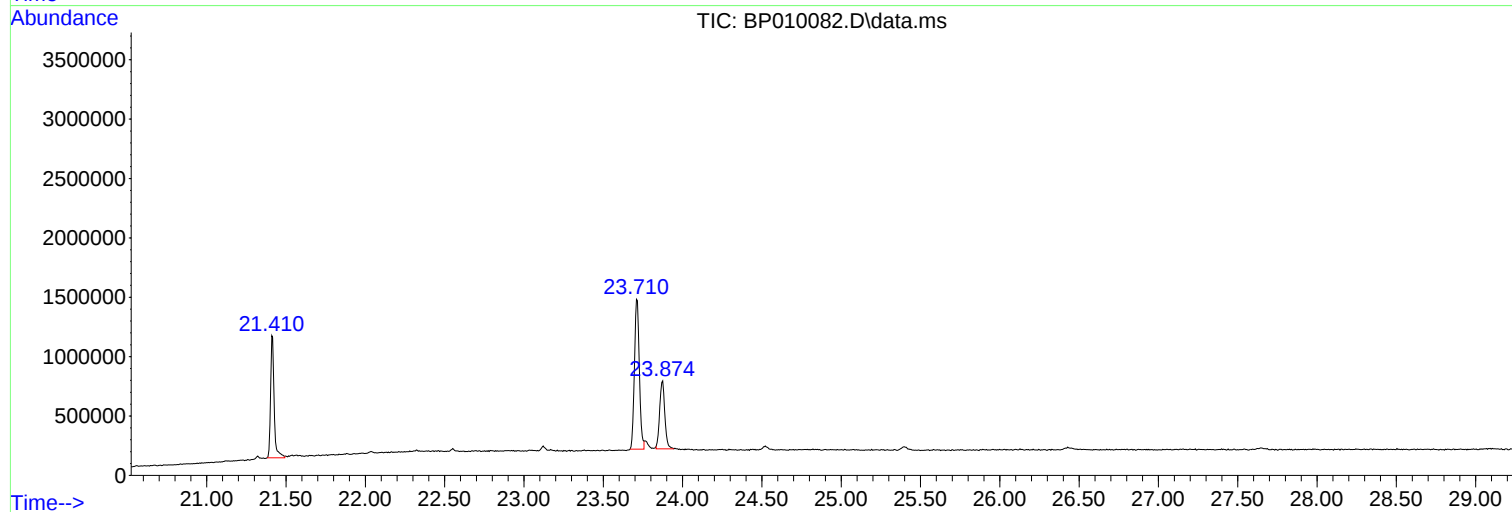
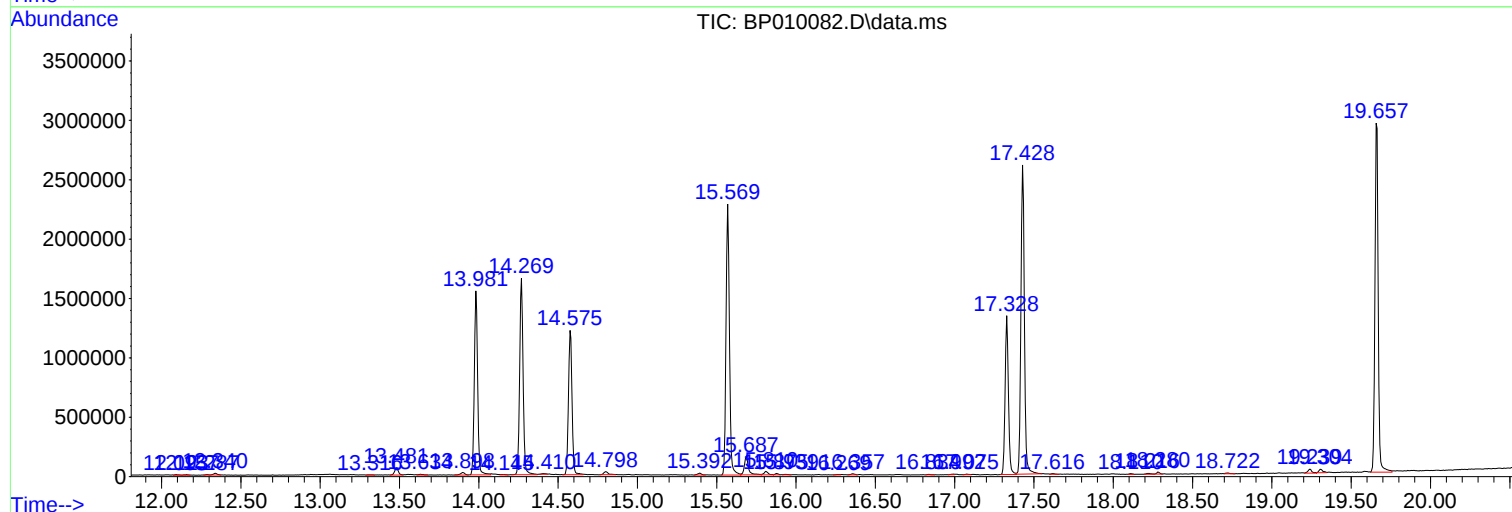
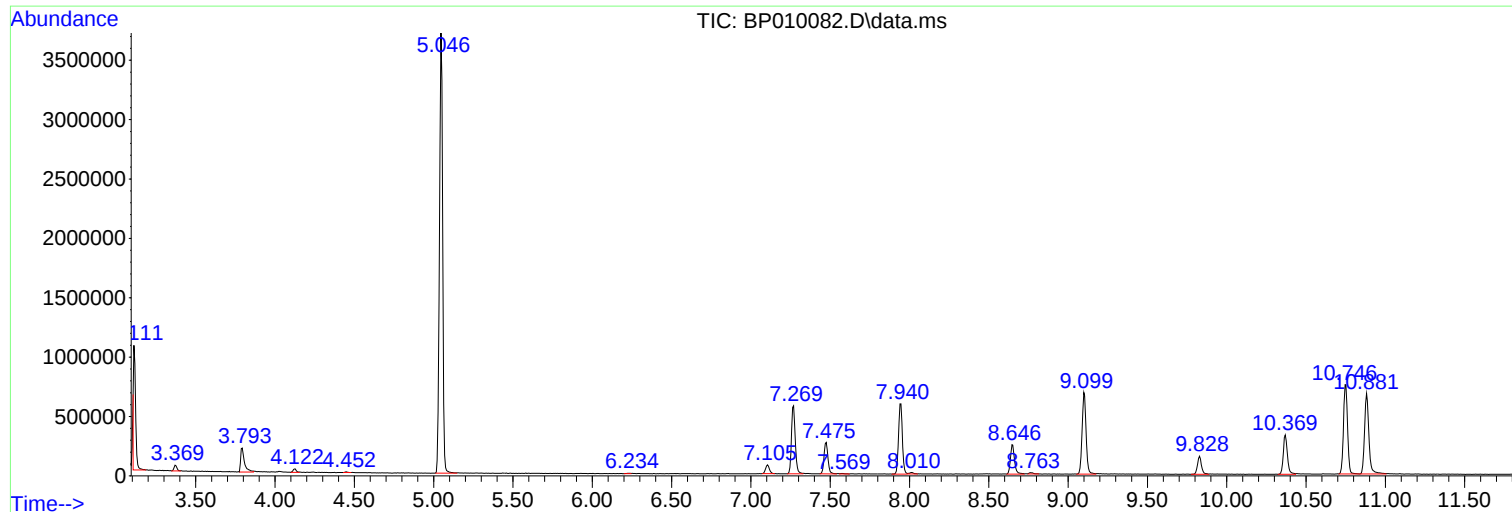
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

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



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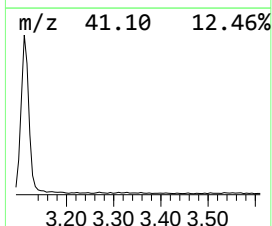
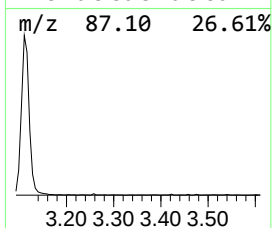
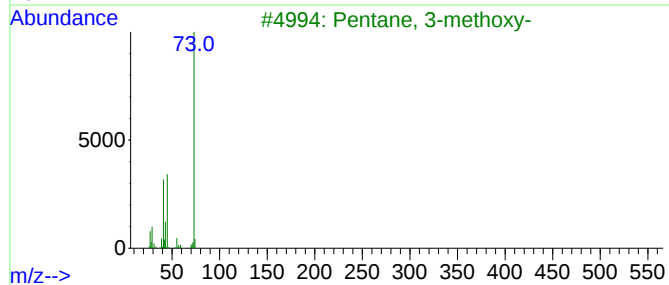
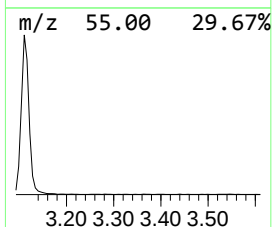
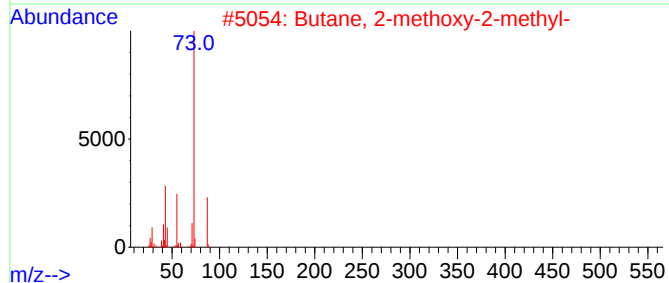
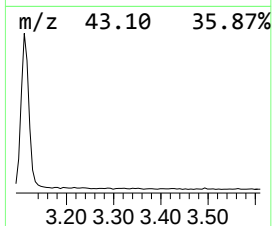
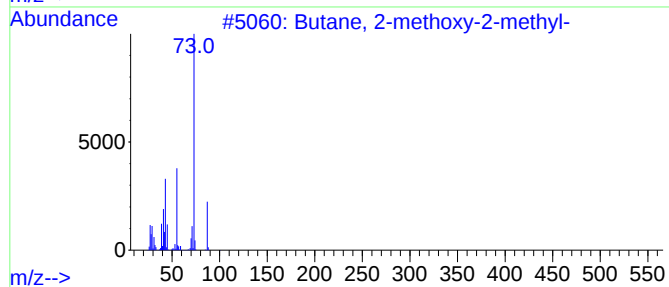
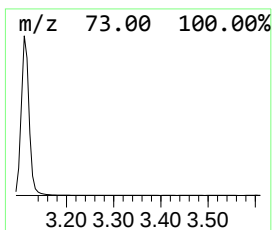
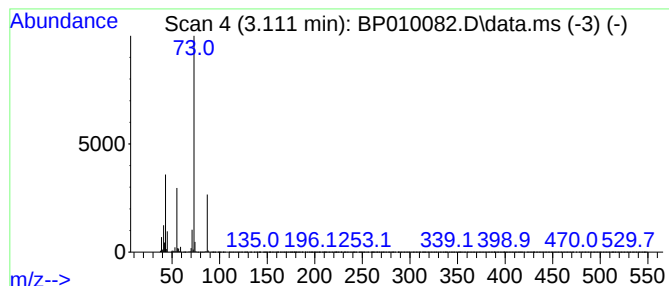
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.111	18.61 ng/ul	915898	1,4-Dichlorobenzene-d4	7.946

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78		
2	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78		
3	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17		
4	Silane, tetramethyl-	88	C4H12Si	000075-76-3	17		
5	N-Ethylformamide	73	C3H7NO	000627-45-2	9		



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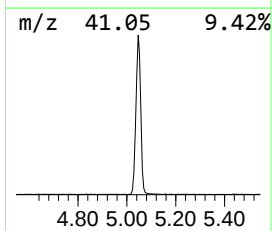
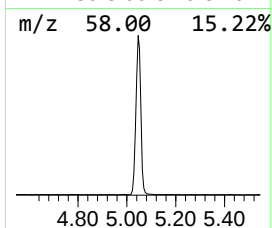
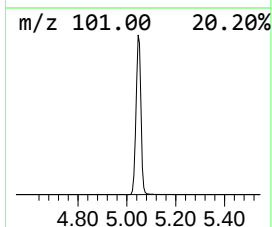
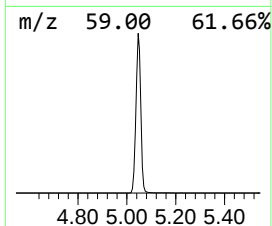
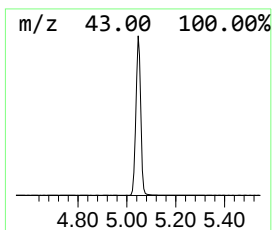
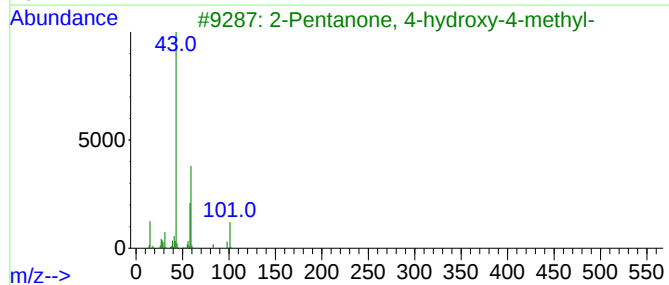
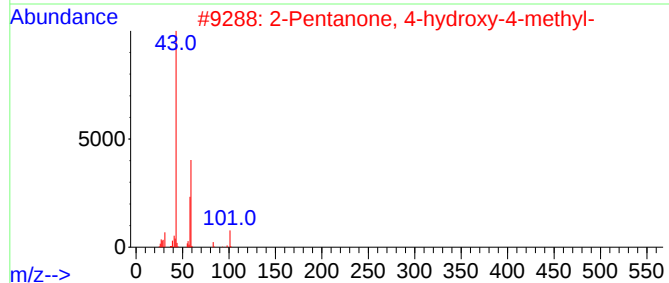
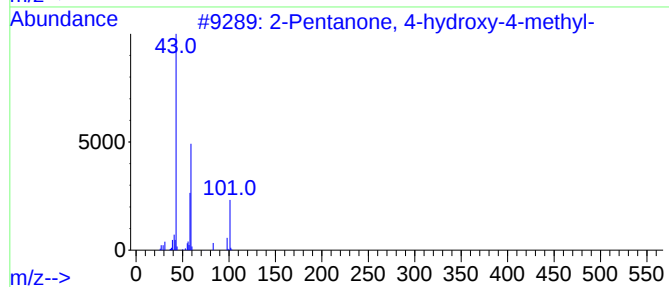
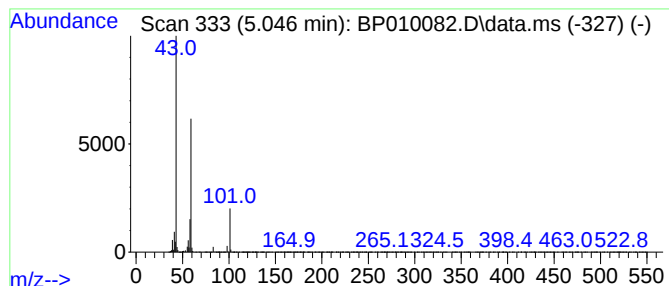
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TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
5.046	103.84 ng/ul	5111990	1,4-Dichlorobenzene-d4	7.946	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
3	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	28
4	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25
5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	23



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-metho...	3.111	18.6	ng/ul	915898	1	7.946	984571	20.0
2-Pentanone, 4-...	5.046	103.8	ng/ul	5111990	1	7.946	984571	20.0