

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041822\
 Data File : BP009923.D
 Acq On : 18 Apr 2022 19:29
 Operator : CG/JU
 Sample : N2421-01
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0J55

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-BP040722.M
 Title : SVOA CALIBRATION

Signal : TIC: BP009923.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.117	3	5	19	rVB	289694	384348	7.88%	0.947%
2	3.381	46	50	55	rBV	20250	27655	0.57%	0.068%
3	3.799	118	121	136	rBV	114930	204138	4.18%	0.503%
4	4.134	174	178	182	rBV4	7335	12334	0.25%	0.030%
5	5.264	362	370	375	rBV2	18612	32613	0.67%	0.080%
6	7.105	677	683	696	rVB	117502	220772	4.52%	0.544%
7	7.275	705	712	722	rBV	414763	667909	13.69%	1.645%
8	7.481	740	747	757	rBV	414866	666326	13.65%	1.641%
9	7.952	819	827	844	rBV	440043	730953	14.98%	1.800%
10	8.205	863	870	874	rBV10	2331	4902	0.10%	0.012%
11	8.305	882	887	893	rBV2	12485	18992	0.39%	0.047%
12	8.652	939	946	960	rBV	365877	591058	12.11%	1.456%
13	8.781	965	968	975	rVB9	3489	5296	0.11%	0.013%
14	9.111	1017	1024	1038	rBV	577128	958416	19.64%	2.360%
15	9.287	1049	1054	1063	rBV	3745	8037	0.16%	0.020%
16	9.569	1097	1102	1109	rVB5	8523	16486	0.34%	0.041%
17	9.834	1138	1147	1162	rBV	457867	762707	15.63%	1.878%
18	10.375	1231	1239	1252	rBV	662467	1107271	22.69%	2.727%
19	10.546	1262	1268	1273	rBV2	23276	41495	0.85%	0.102%
20	10.622	1276	1281	1288	rVB3	16839	30735	0.63%	0.076%
21	10.757	1296	1304	1319	rBV	660265	1126980	23.09%	2.775%
22	10.887	1319	1326	1339	rBV	622774	1084010	22.21%	2.670%
23	12.287	1560	1564	1568	rVB7	3269	5030	0.10%	0.012%
24	12.340	1568	1573	1580	rVB3	16412	26463	0.54%	0.065%
25	12.557	1605	1610	1612	rBV6	3934	5341	0.11%	0.013%
26	12.957	1674	1678	1682	rVB6	4332	6618	0.14%	0.016%
27	13.199	1716	1719	1725	rVB7	7585	10618	0.22%	0.026%
28	13.393	1744	1752	1756	rBV4	21627	38739	0.79%	0.095%
29	13.557	1775	1780	1782	rBV6	5690	8866	0.18%	0.022%
30	13.587	1782	1785	1791	rVB2	9196	14355	0.29%	0.035%
31	13.993	1847	1854	1860	rBV	1576374	2232080	45.74%	5.497%
32	14.234	1890	1895	1896	rBV2	26417	32318	0.66%	0.080%
33	14.275	1896	1902	1917	rVB	1572500	2353766	48.23%	5.797%
34	14.493	1935	1939	1944	rVB6	5992	9144	0.19%	0.023%
35	14.581	1947	1954	1974	rVB	1104003	1712293	35.09%	4.217%
36	14.763	1976	1985	1995	rBV	126981	211393	4.33%	0.521%

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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_P\Methods\SFAM-EPA-BP040722.M
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37	15.222	2058	2063	2069	rVB2	12111	19812	0.41%	0.049%
38	15.393	2087	2092	2097	rBV2	31168	48748	1.00%	0.120%
39	15.440	2097	2100	2105	rVB5	7987	11930	0.24%	0.029%
40	15.575	2116	2123	2135	rBV	2384412	3383745	69.34%	8.333%
41	15.687	2135	2142	2158	rVV	821626	1154034	23.65%	2.842%
42	15.816	2158	2164	2171	rVV2	26270	40014	0.82%	0.099%
43	15.940	2180	2185	2190	rBV7	9237	13490	0.28%	0.033%
44	16.134	2212	2218	2225	rBV7	5302	12373	0.25%	0.030%
45	16.222	2228	2233	2236	rBV6	5359	9421	0.19%	0.023%
46	16.263	2236	2240	2244	rVV5	13658	23002	0.47%	0.057%
47	16.304	2244	2247	2250	rVV4	11219	15718	0.32%	0.039%
48	16.363	2254	2257	2264	rVB6	11751	19635	0.40%	0.048%
49	16.481	2273	2277	2282	rVB4	10116	13390	0.27%	0.033%
50	16.651	2301	2306	2307	rBV5	3860	5217	0.11%	0.013%
51	16.816	2330	2334	2336	rBV5	7444	8707	0.18%	0.021%
52	17.016	2361	2368	2372	rBV9	6007	13862	0.28%	0.034%
53	17.104	2377	2383	2389	rBV6	14827	30451	0.62%	0.075%
54	17.263	2405	2410	2415	rVB4	27627	37023	0.76%	0.091%
55	17.334	2415	2422	2432	rBV	1491780	2184601	44.77%	5.380%
56	17.434	2432	2439	2449	rVV	2801411	4118189	84.39%	10.142%
57	17.528	2453	2455	2460	rVB6	7902	9432	0.19%	0.023%
58	17.651	2468	2476	2478	rBV8	9123	20080	0.41%	0.049%
59	17.845	2506	2509	2512	rBV5	6572	9058	0.19%	0.022%
60	17.963	2523	2529	2531	rVB7	3586	6184	0.13%	0.015%
61	18.010	2531	2537	2542	rBV4	16542	25372	0.52%	0.062%
62	18.116	2553	2555	2558	rBV4	4777	6957	0.14%	0.017%
63	18.216	2568	2572	2579	rBV3	45181	66051	1.35%	0.163%
64	18.287	2581	2584	2589	rVB3	13830	19072	0.39%	0.047%
65	18.516	2619	2623	2626	rBV5	7764	14774	0.30%	0.036%
66	18.751	2661	2663	2667	rVB5	8034	7676	0.16%	0.019%
67	18.969	2696	2700	2703	rBV5	8338	11525	0.24%	0.028%
68	19.051	2711	2714	2717	rBV3	11164	16109	0.33%	0.040%
69	19.239	2742	2746	2750	rBV2	42192	61504	1.26%	0.151%
70	19.281	2750	2753	2754	rVV3	14040	15421	0.32%	0.038%
71	19.304	2754	2757	2764	rVB	35899	52233	1.07%	0.129%
72	19.445	2779	2781	2786	rBV3	10058	10781	0.22%	0.027%
73	19.663	2812	2818	2828	rBV	3736390	4879826	100.00%	12.017%
74	21.122	3062	3066	3074	rBV	156906	225364	4.62%	0.555%
75	21.416	3111	3116	3124	rBV	1826354	2339147	47.94%	5.761%
76	23.722	3500	3508	3515	rBV	2139303	4154744	85.14%	10.232%

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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

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Title : SVOA CALIBRATION

77	23.874	3527	3534	3545	rVB2	1016314	2120891	43.46%	5.223%
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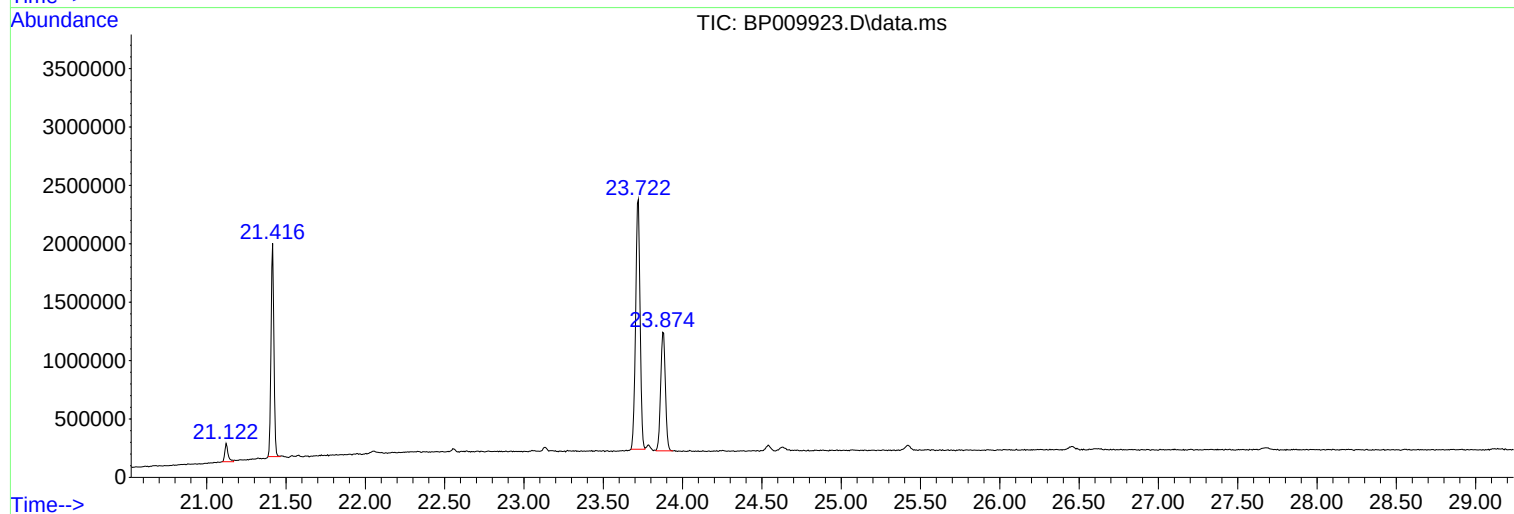
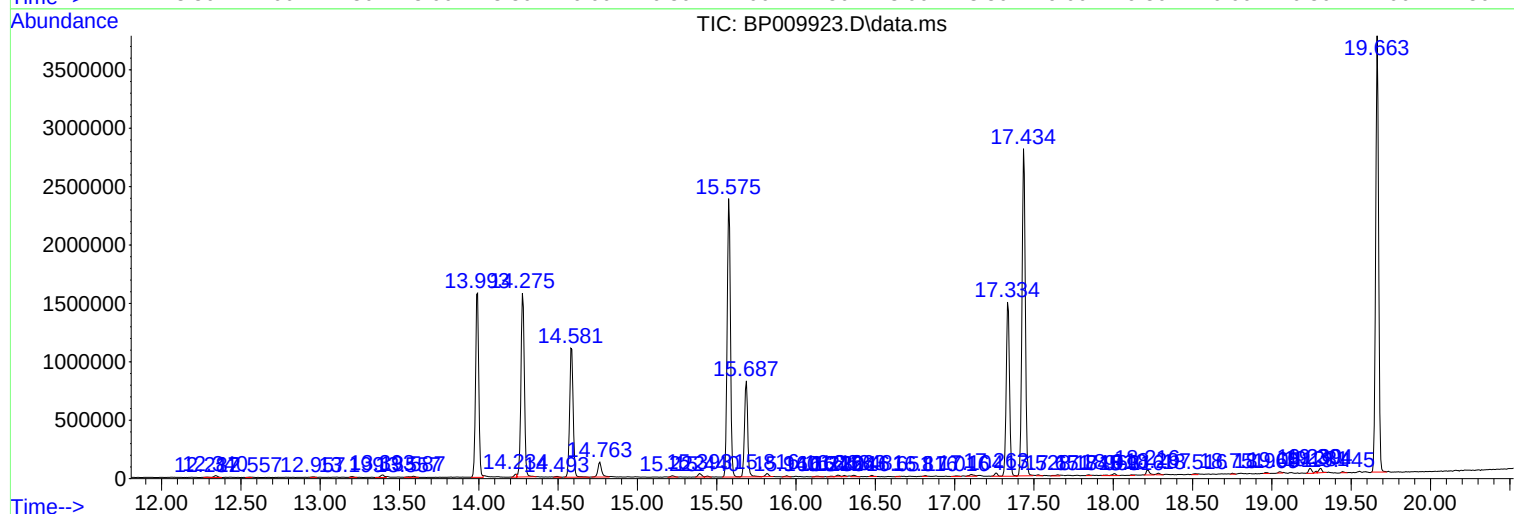
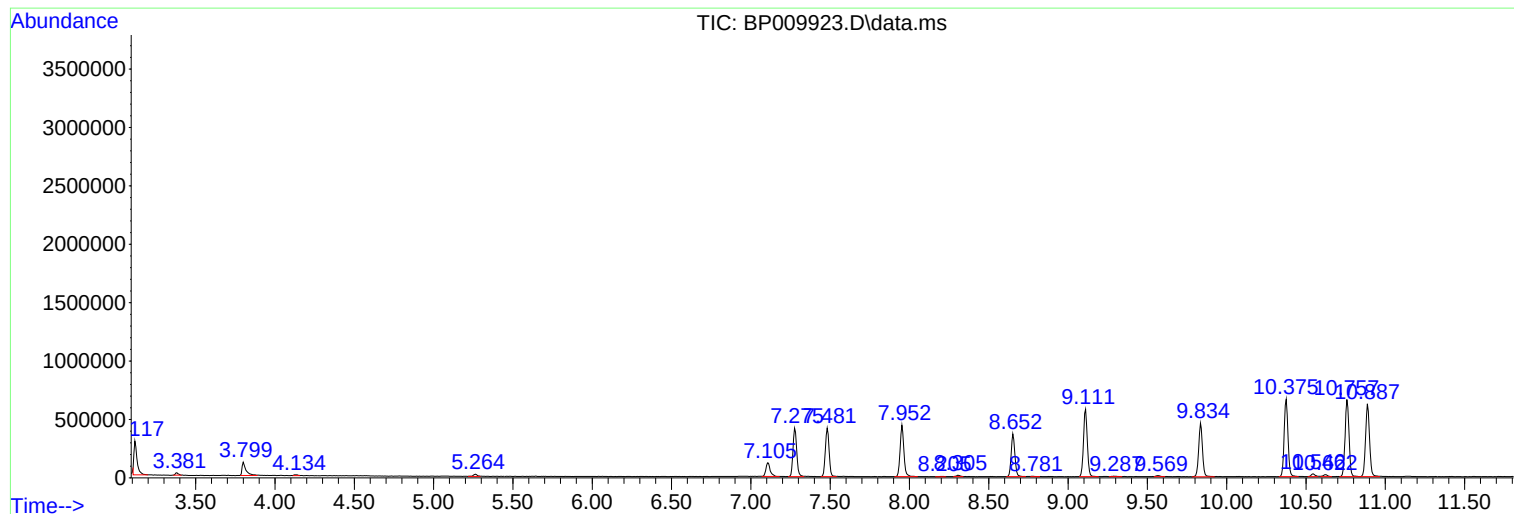
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION

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



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Instrument :
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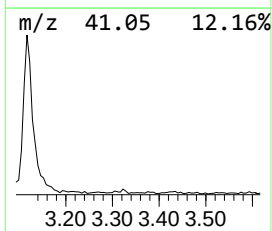
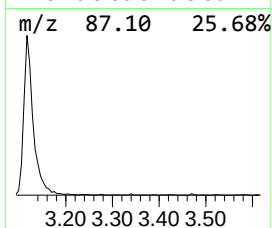
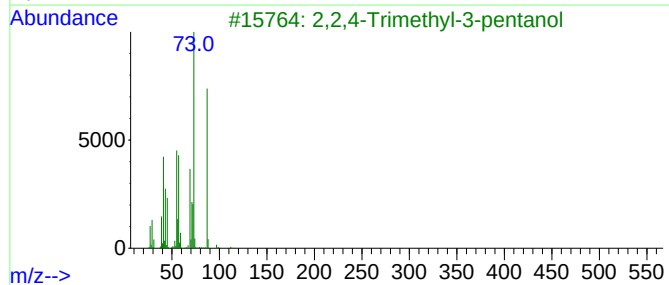
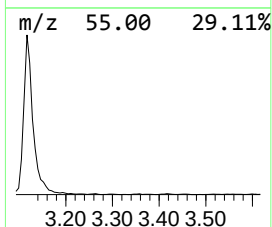
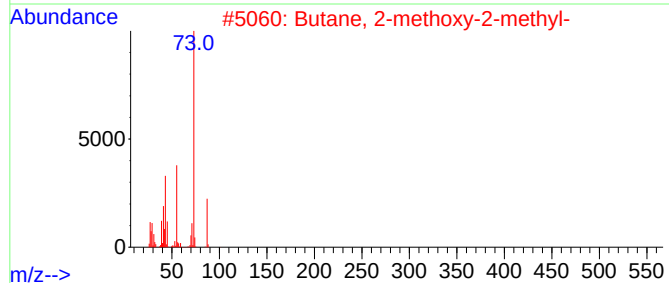
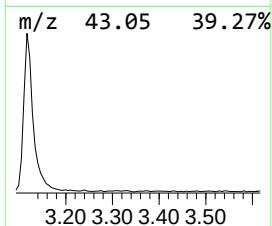
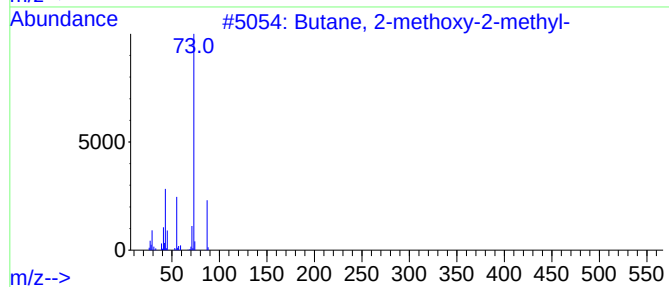
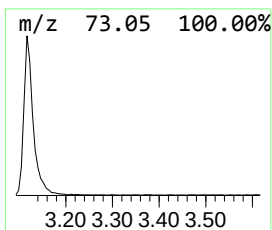
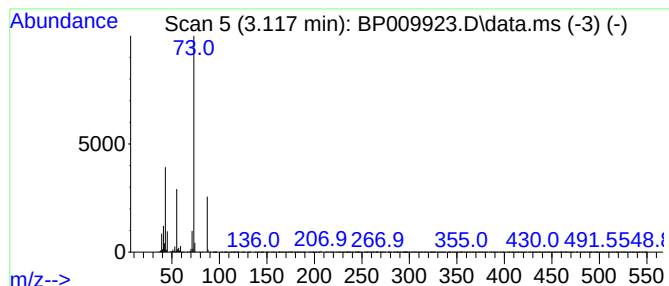
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION

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.
3.117	10.52 ng/ul	384348	1,4-Dichlorobenzene-d4	7.952

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	40
4			Silane, tetramethyl-	88	C4H12Si	000075-76-3	38
5			Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	33



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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...	3.117	10.5	ng/ul	384348	1	7.952	730953	20.0