

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100322\
 Data File : BP011872.D
 Acq On : 03 Oct 2022 15:07
 Operator : CG/JU
 Sample : N4749-07
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 CB8Q1

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-BP093022.M
 Title : SVOA CALIBRATION

Signal : TIC: BP011872.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.034	4	8	30	rVB	5167652	6806899	91.66%	10.141%
2	3.305	50	54	59	rBV	59711	69992	0.94%	0.104%
3	3.734	124	127	141	rBV	225174	353573	4.76%	0.527%
4	3.834	141	144	149	rVB3	14656	22051	0.30%	0.033%
5	3.958	160	165	173	rVB	33353	53716	0.72%	0.080%
6	4.052	177	181	185	rBV	17135	24291	0.33%	0.036%
7	4.281	216	220	226	rVB2	13547	19310	0.26%	0.029%
8	4.505	255	258	264	rVB7	5048	8085	0.11%	0.012%
9	4.693	286	290	295	rBV4	6464	11912	0.16%	0.018%
10	4.993	335	341	348	rVB3	10776	21560	0.29%	0.032%
11	5.164	366	370	376	rVB6	4105	7738	0.10%	0.012%
12	7.058	684	692	705	rBV	237923	443856	5.98%	0.661%
13	7.222	713	720	736	rBV	889406	1447251	19.49%	2.156%
14	7.422	746	754	765	rBV	847397	1390775	18.73%	2.072%
15	7.893	823	834	842	rBV	917496	1524418	20.53%	2.271%
16	7.952	842	844	849	rVB3	4453	7874	0.11%	0.012%
17	8.605	946	955	968	rBV	652843	1150410	15.49%	1.714%
18	9.057	1025	1032	1048	rBV	916952	1566253	21.09%	2.333%
19	9.469	1096	1102	1107	rBV	18791	33216	0.45%	0.049%
20	9.781	1147	1155	1168	rBV	708455	1187922	16.00%	1.770%
21	10.322	1238	1247	1263	rBV	1068944	1929184	25.98%	2.874%
22	10.493	1271	1276	1291	rVB6	15090	43916	0.59%	0.065%
23	10.699	1302	1311	1327	rBV	1251576	2161766	29.11%	3.221%
24	10.840	1327	1335	1356	rVV	610493	1176049	15.84%	1.752%
25	10.987	1358	1360	1371	rVB	6116	12478	0.17%	0.019%
26	11.522	1445	1451	1455	rBV7	10743	19838	0.27%	0.030%
27	12.116	1545	1552	1558	rBV9	6430	11910	0.16%	0.018%
28	12.293	1573	1582	1590	rBV	22064	41987	0.57%	0.063%
29	12.493	1612	1616	1621	rBV7	3580	7474	0.10%	0.011%
30	13.263	1742	1747	1752	rBV8	3787	8299	0.11%	0.012%
31	13.357	1759	1763	1767	rBV	17354	25773	0.35%	0.038%
32	13.434	1771	1776	1781	rVV	28742	44991	0.61%	0.067%
33	13.510	1785	1789	1797	rVB10	4844	9142	0.12%	0.014%
34	13.716	1817	1824	1829	rBV2	10440	16870	0.23%	0.025%
35	13.945	1856	1863	1876	rBV	2147834	3008732	40.52%	4.482%
36	14.228	1899	1911	1925	rBV	2477984	3732296	50.26%	5.560%

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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_P\Methods\SFAM-EPA-BP093022.M
 Title : SVOA CALIBRATION

37	14.428	1943	1945	1953	rVB9	8516	13249	0.18%	0.020%
38	14.534	1955	1963	1969	rBV	1730060	2578624	34.72%	3.842%
39	14.587	1969	1972	1987	rVB	53866	131044	1.76%	0.195%
40	14.745	1993	1999	2020	rBV	115925	282545	3.80%	0.421%
41	14.898	2021	2025	2031	rVV7	10646	20585	0.28%	0.031%
42	14.951	2031	2034	2040	rVB6	8589	14756	0.20%	0.022%
43	15.287	2087	2091	2094	rVB	11488	14217	0.19%	0.021%
44	15.357	2098	2103	2111	rVB	164416	224398	3.02%	0.334%
45	15.534	2126	2133	2138	rBV	3108329	4779098	64.36%	7.120%
46	15.645	2146	2152	2162	rBV	855632	1173806	15.81%	1.749%
47	15.769	2169	2173	2180	rVB4	17479	29261	0.39%	0.044%
48	17.051	2388	2391	2395	rBV5	10987	11933	0.16%	0.018%
49	17.292	2425	2432	2437	rBV	1947513	2788772	37.55%	4.155%
50	17.334	2437	2439	2443	rVV	101099	113734	1.53%	0.169%
51	17.392	2443	2449	2459	rVB	3650277	5199006	70.01%	7.745%
52	17.510	2466	2469	2476	rVB8	17100	27639	0.37%	0.041%
53	17.657	2490	2494	2502	rBV5	35917	58978	0.79%	0.088%
54	18.175	2577	2582	2588	rBV2	93821	145691	1.96%	0.217%
55	19.204	2753	2757	2761	rVB3	48290	64206	0.86%	0.096%
56	19.269	2765	2768	2775	rBV	56716	89248	1.20%	0.133%
57	19.404	2788	2791	2801	rBV5	50546	76559	1.03%	0.114%
58	19.628	2823	2829	2841	rBV	4642593	6453007	86.90%	9.614%
59	21.080	3073	3076	3082	rBV3	200450	256151	3.45%	0.382%
60	21.374	3122	3126	3132	rVB	2546032	3185971	42.90%	4.746%
61	23.651	3507	3513	3527	rVB2	3574969	7425968	100.00%	11.063%
62	23.810	3534	3540	3548	rVB2	1792169	3563715	47.99%	5.309%

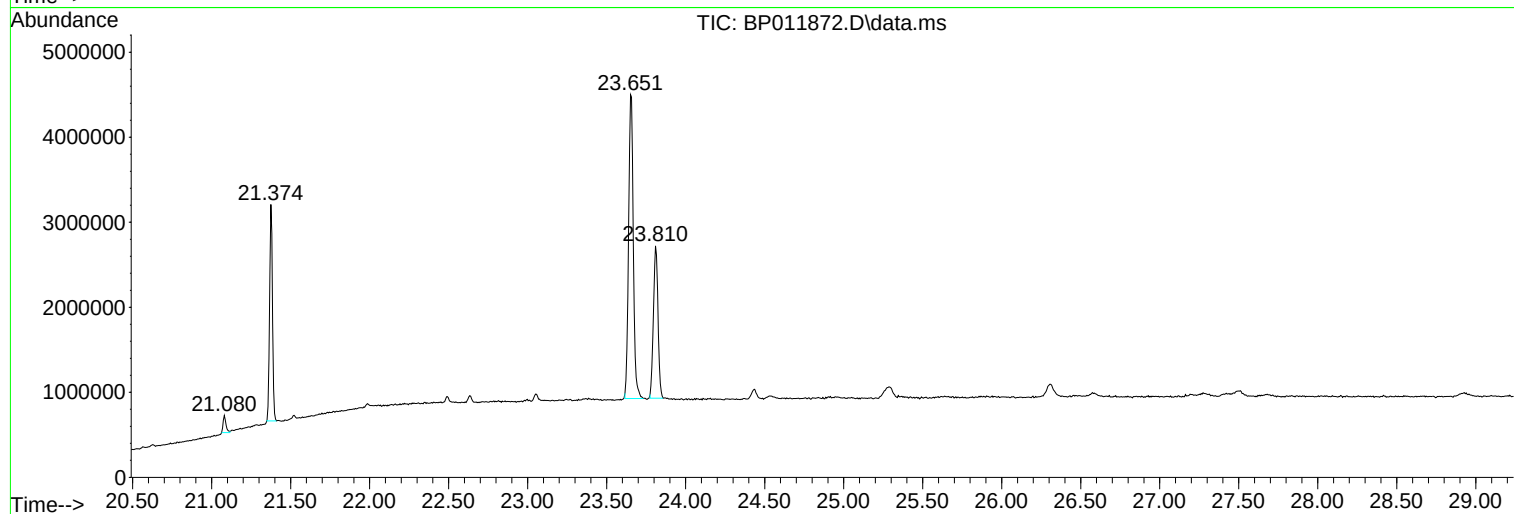
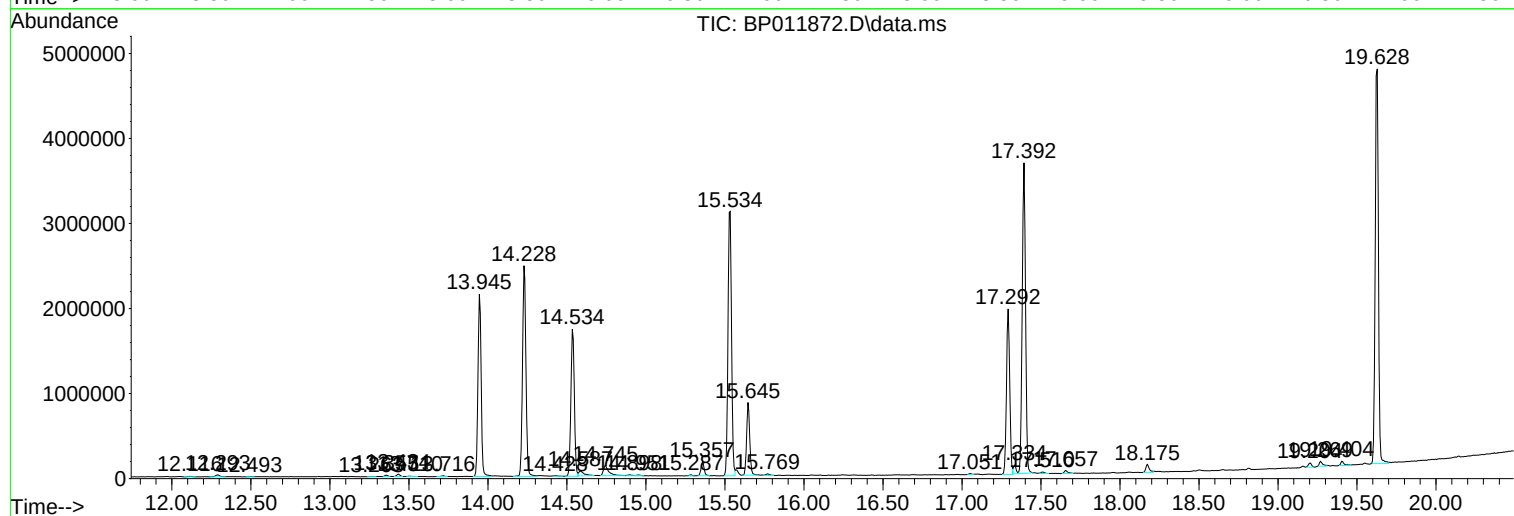
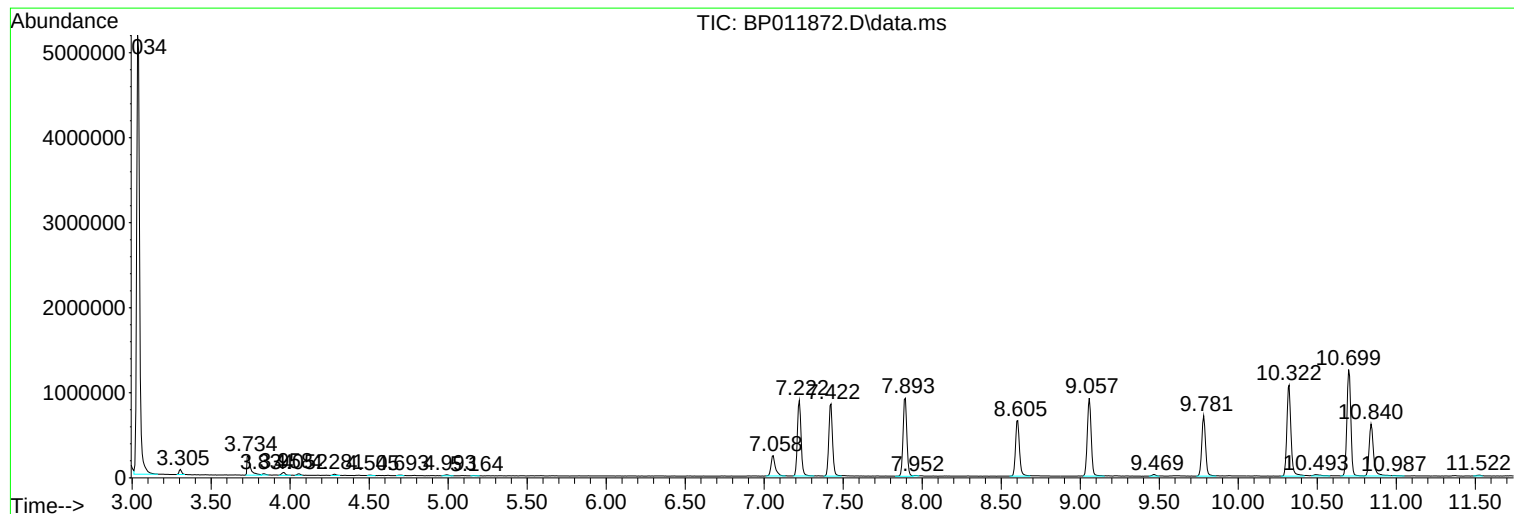
Sum of corrected areas: 67123968

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

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



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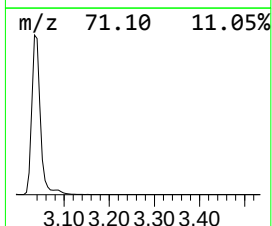
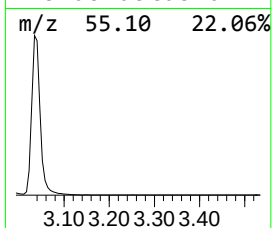
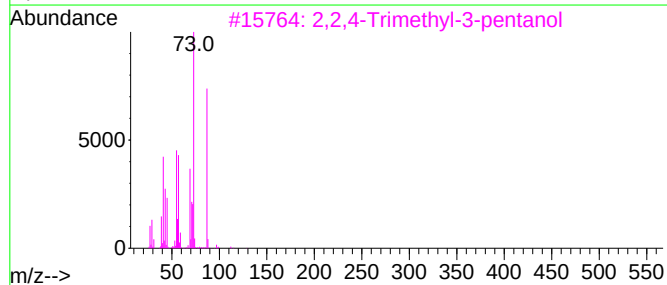
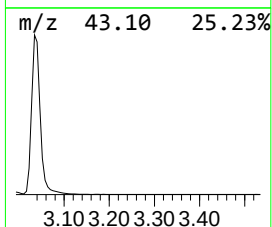
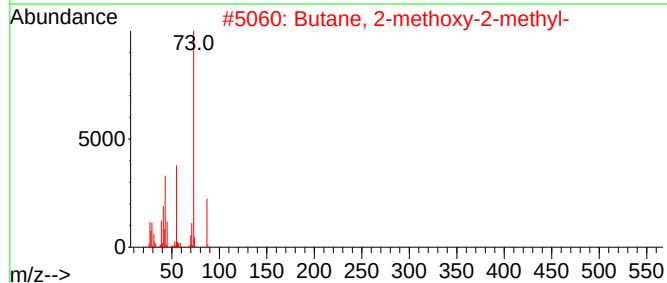
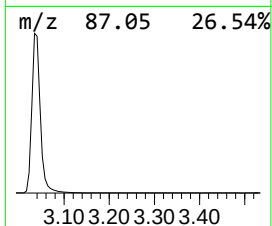
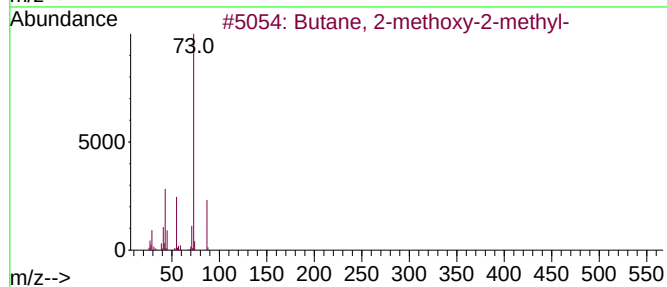
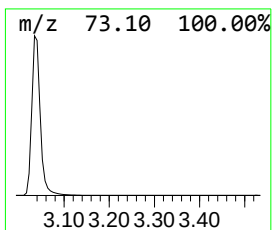
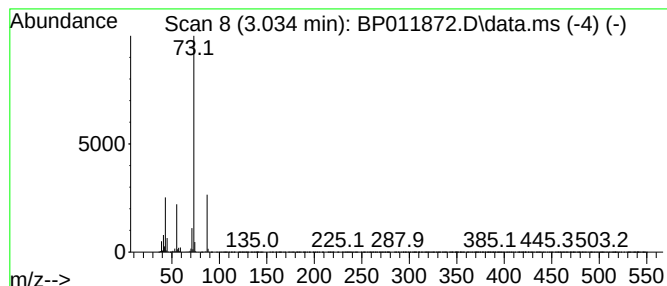
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST0.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.
3.034	89.30 ng/ul	6806900	1,4-Dichlorobenzene-d4	7.893

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	86
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			N-Ethylformamide	73	C3H7NO	000627-45-2	9
5			Silane, tetramethyl-	88	C4H12Si	000075-76-3	9



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TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-metho...	3.034	89.3	ng/ul	6806900	1	7.893	1524420	20.0