

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101821\
 Data File : BP007469.D
 Acq On : 18 Oct 2021 18:04
 Operator : CG/JU
 Sample : PB139924BL
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK924

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-BP101421.M
 Title : SVOA CALIBRATION

Signal : TIC: BP007469.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	18	rVB	188947	248151	4.81%	0.495%
2	3.381	46	50	58	rVB	56888	78754	1.53%	0.157%
3	3.793	116	120	134	rBV	691690	1005467	19.50%	2.008%
4	4.134	174	178	182	rVB	13180	16239	0.32%	0.032%
5	7.116	678	685	694	rBV	776796	1239894	24.05%	2.476%
6	7.281	707	713	724	rVB	593765	967865	18.78%	1.932%
7	7.487	741	748	758	rBV	805652	1251053	24.27%	2.498%
8	7.958	821	828	837	rVB	647731	1020617	19.80%	2.038%
9	8.663	939	948	955	rBV	961179	1565290	30.36%	3.125%
10	9.110	1015	1024	1032	rBV	726387	1194583	23.17%	2.385%
11	9.840	1140	1148	1156	rBV	554039	912186	17.70%	1.821%
12	10.381	1232	1240	1252	rBV	957984	1569689	30.45%	3.134%
13	10.763	1296	1305	1316	rVB	877594	1463689	28.39%	2.922%
14	10.893	1318	1327	1340	rBV	925502	1563130	30.32%	3.121%
15	12.346	1568	1574	1582	rBV2	19395	34543	0.67%	0.069%
16	13.569	1776	1782	1786	rVB8	3697	5576	0.11%	0.011%
17	14.004	1849	1856	1871	rBV	1670796	2370960	45.99%	4.734%
18	14.287	1896	1904	1911	rBV	1840583	2849132	55.27%	5.689%
19	14.592	1949	1956	1965	rBV	1341844	2000054	38.80%	3.993%
20	14.775	1980	1987	1999	rBV	854090	1240809	24.07%	2.477%
21	15.587	2118	2125	2134	rBV	2471513	3520673	68.30%	7.029%
22	15.698	2137	2144	2155	rBV	622800	866166	16.80%	1.729%
23	15.828	2160	2166	2172	rVB2	30047	48073	0.93%	0.096%
24	15.951	2182	2187	2190	rBV7	4733	6096	0.12%	0.012%
25	16.269	2235	2241	2243	rBV7	5128	9228	0.18%	0.018%
26	16.375	2256	2259	2264	rVB7	4586	7445	0.14%	0.015%
27	17.028	2367	2370	2376	rVB6	4247	5853	0.11%	0.012%
28	17.134	2382	2388	2397	rVB3	8591	18366	0.36%	0.037%
29	17.345	2418	2424	2432	rBV	1724162	2480843	48.13%	4.953%
30	17.445	2433	2441	2452	rBV	2905793	4119768	79.92%	8.226%
31	17.581	2460	2464	2471	rVB9	7505	14989	0.29%	0.030%
32	17.716	2483	2487	2491	rBV4	5701	8168	0.16%	0.016%
33	18.022	2536	2539	2544	rVB7	4997	5827	0.11%	0.012%
34	18.239	2572	2576	2582	rBV9	8194	13625	0.26%	0.027%
35	19.257	2745	2749	2754	rVB2	44486	58490	1.13%	0.117%
36	19.322	2756	2760	2767	rBV	44808	67260	1.30%	0.134%

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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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37	19.680	2815	2821	2829	rBV	3684299	4959551	96.21%	9.902%
38	21.433	3114	3119	3126	rBV	2348601	2987187	57.95%	5.964%
39	23.733	3502	3510	3519	rBV2	2570981	5154938	100.00%	10.292%
40	23.892	3530	3537	3546	rVB2	1473201	3134605	60.81%	6.259%

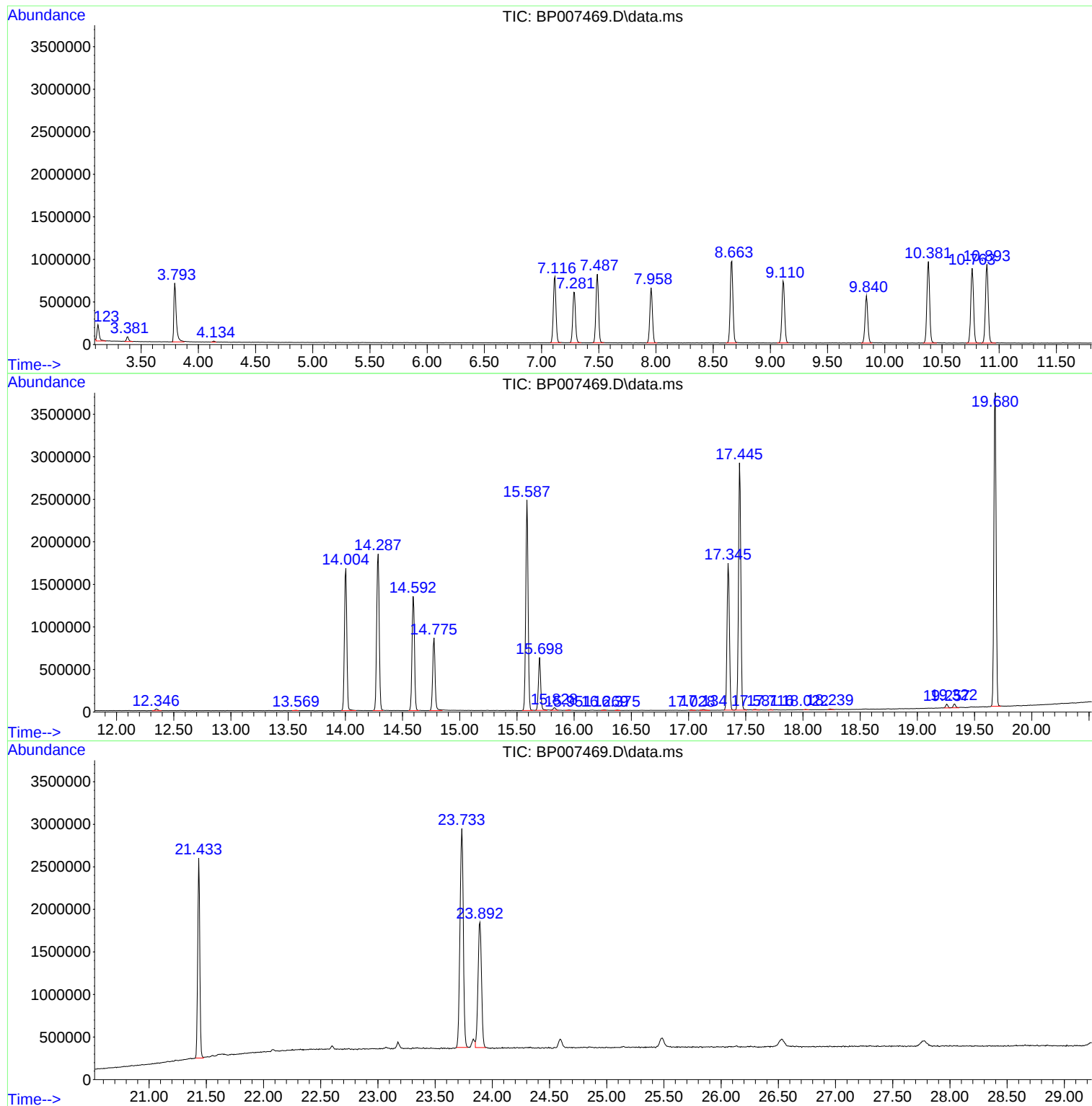
Sum of corrected areas: 50084832

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP101421.M
Quant Title : SVOA CALIBRATION

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



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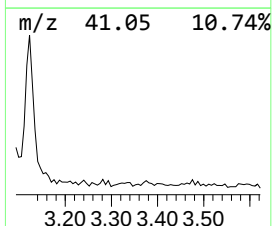
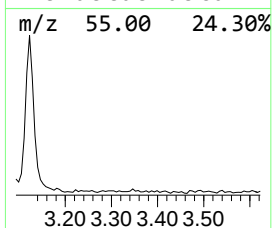
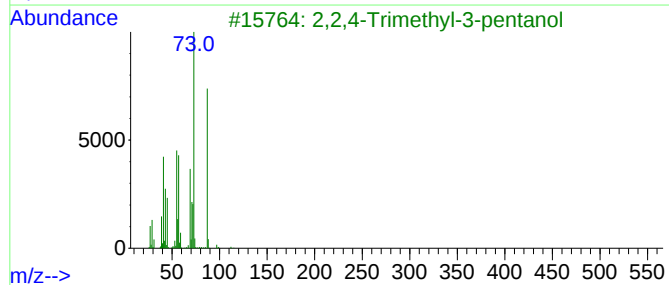
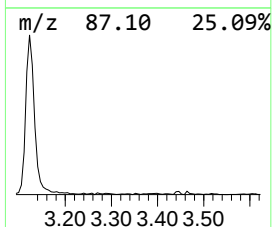
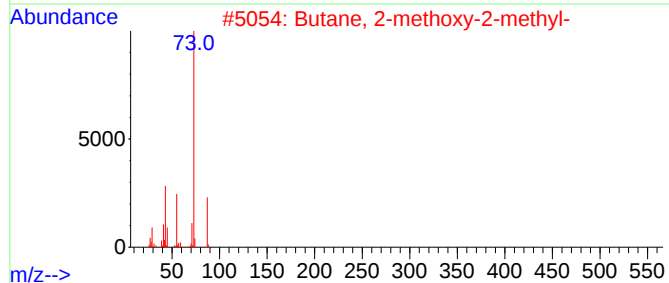
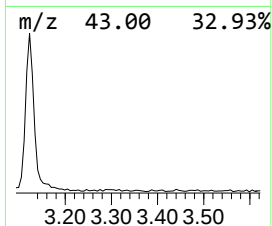
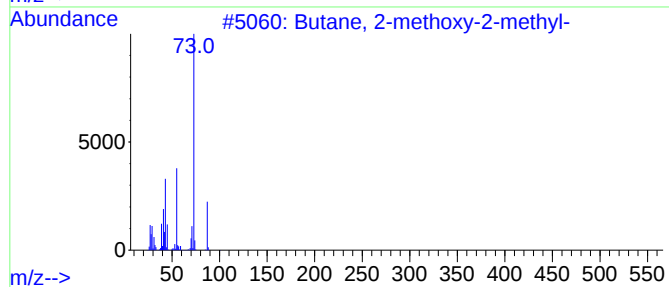
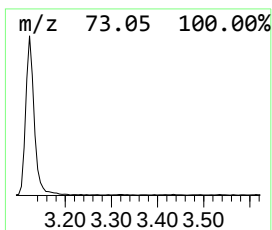
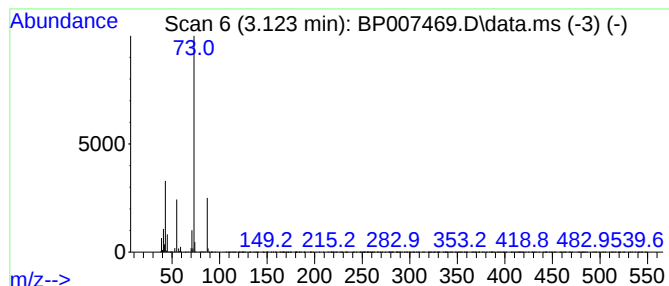
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP101421.M
Quant Title : SVOA CALIBRATION

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	4.86 ng/ul	248151	1,4-Dichlorobenzene-d4	7.958

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	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40		
4	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17		
5	N-Ethylformamide	73	C3H7NO	000627-45-2	9		



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Quant Title : SVOA CALIBRATION

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	4.9	ng/u1	248151	1	7.958	1020620	20.0