

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041522\
 Data File : BP009862.D
 Acq On : 15 Apr 2022 15:37
 Operator : CG/JU
 Sample : N2422-02
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0J57

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: BP009862.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.122	3	6	22	rVB	307006	465123	10.37%	1.214%
2	3.381	46	50	56	rBV	23334	33632	0.75%	0.088%
3	3.805	118	122	139	rBV	92273	174540	3.89%	0.456%
4	4.134	175	178	183	rVB3	6077	9150	0.20%	0.024%
5	7.110	676	684	694	rBV	136453	218820	4.88%	0.571%
6	7.281	707	713	725	rBV	416160	672331	15.00%	1.755%
7	7.487	740	748	762	rBV	429282	696088	15.53%	1.817%
8	7.958	820	828	838	rBV	434022	704970	15.72%	1.840%
9	8.657	940	947	961	rBV	371949	617998	13.78%	1.613%
10	9.116	1014	1025	1037	rBV	557402	951310	21.22%	2.483%
11	9.575	1097	1103	1108	rVB4	8243	14287	0.32%	0.037%
12	9.840	1138	1148	1159	rBV	431023	731134	16.31%	1.908%
13	10.381	1233	1240	1254	rBV	613668	1049297	23.41%	2.739%
14	10.628	1277	1282	1289	rVB	3152	7508	0.17%	0.020%
15	10.763	1295	1305	1319	rBV	611149	1072603	23.93%	2.800%
16	10.893	1319	1327	1341	rVV	644753	1108599	24.73%	2.893%
17	12.198	1546	1549	1555	rVB8	2614	4836	0.11%	0.013%
18	12.351	1570	1575	1580	rVB	13254	22061	0.49%	0.058%
19	13.210	1717	1721	1727	rVB9	3106	5522	0.12%	0.014%
20	13.398	1746	1753	1757	rBV5	13322	23991	0.54%	0.063%
21	13.428	1757	1758	1763	rVB4	6592	7006	0.16%	0.018%
22	13.992	1847	1854	1867	rBV	1401884	2079206	46.38%	5.427%
23	14.281	1896	1903	1920	rVB	1525904	2196171	48.99%	5.732%
24	14.587	1947	1955	1966	rBV	1027203	1547057	34.51%	4.038%
25	14.763	1978	1985	1994	rBV	149750	248108	5.53%	0.648%
26	15.398	2088	2093	2099	rBV2	11811	17639	0.39%	0.046%
27	15.581	2116	2124	2135	rBV	2072817	3057266	68.19%	7.980%
28	15.692	2136	2143	2157	rVV	728702	1008449	22.49%	2.632%
29	15.822	2160	2165	2172	rVB3	24747	39507	0.88%	0.103%
30	16.269	2238	2241	2244	rBV5	3606	4880	0.11%	0.013%
31	16.369	2254	2258	2263	rVB4	8976	11976	0.27%	0.031%
32	17.016	2365	2368	2374	rBV7	4357	6135	0.14%	0.016%
33	17.128	2382	2387	2390	rBV6	3340	5953	0.13%	0.016%
34	17.339	2416	2423	2432	rBV	1363611	1964651	43.82%	5.128%
35	17.439	2432	2440	2451	rBV	2638331	3693264	82.38%	9.640%
36	18.010	2535	2537	2541	rBV5	4754	5617	0.13%	0.015%

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37	18.222	2569	2573	2580	rBV2	49906	73592	1.64%	0.192%
38	18.286	2582	2584	2589	rVB6	9334	12990	0.29%	0.034%
39	19.063	2711	2716	2718	rBV6	5002	9200	0.21%	0.024%
40	19.245	2743	2747	2752	rBV3	37448	49379	1.10%	0.129%
41	19.316	2754	2759	2765	rBV	38068	52870	1.18%	0.138%
42	19.457	2779	2783	2786	rBV6	11828	18037	0.40%	0.047%
43	19.604	2804	2808	2811	rBV6	9532	13283	0.30%	0.035%
44	19.669	2813	2819	2831	rVB	3479920	4483182	100.00%	11.701%
45	21.127	3063	3067	3074	rBV2	123665	180960	4.04%	0.472%
46	21.422	3112	3117	3124	rBV	1795247	2290563	51.09%	5.978%
47	23.727	3501	3509	3517	rBV2	2291383	4461277	99.51%	11.644%
48	23.886	3529	3536	3546	rVB2	1080447	2191735	48.89%	5.720%

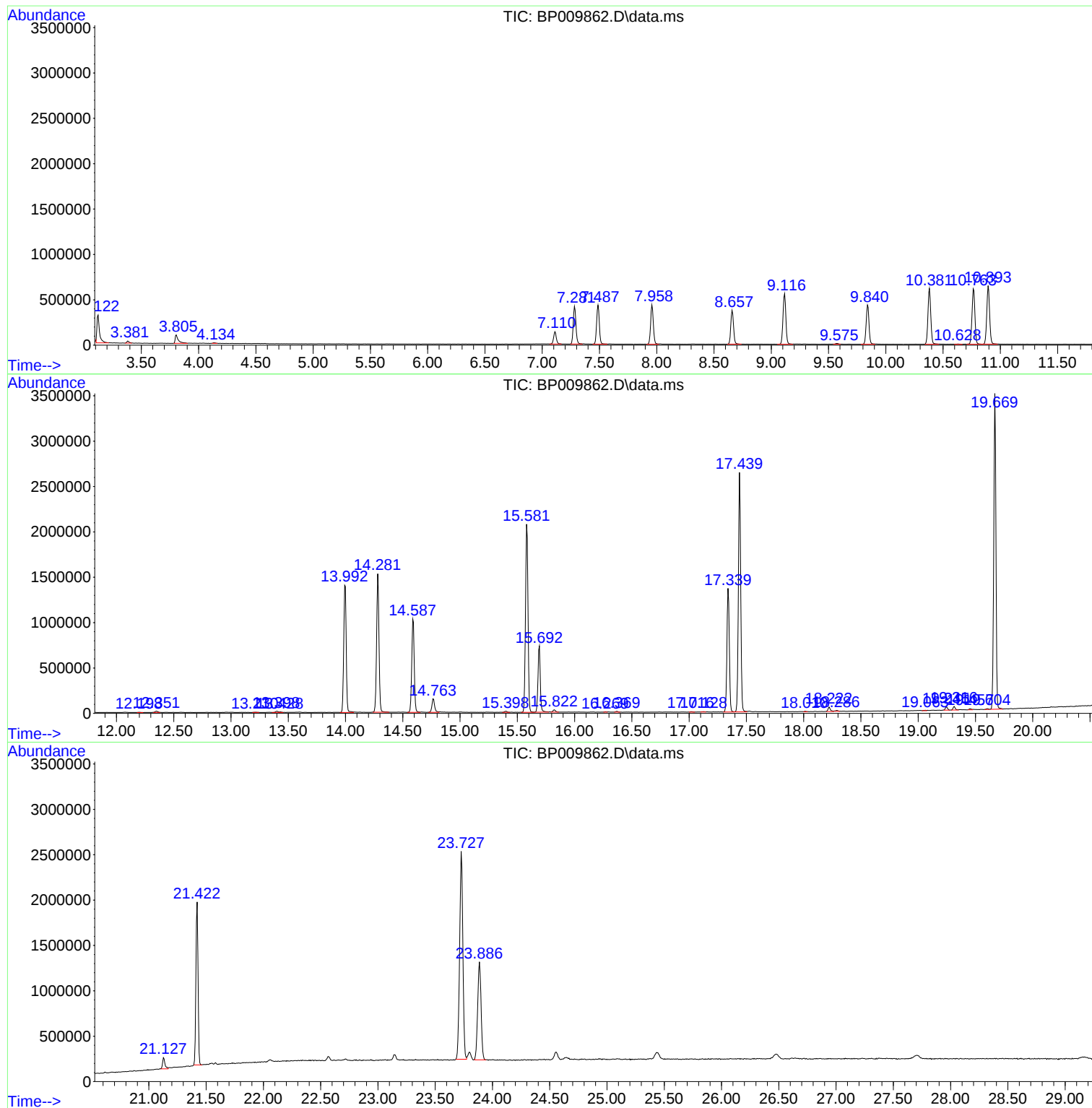
Sum of corrected areas: 38313753

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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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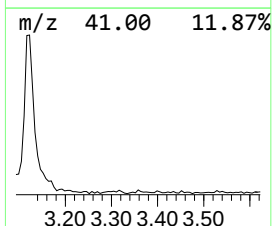
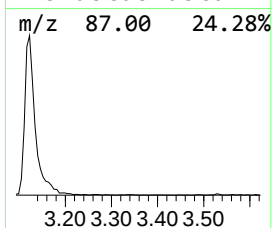
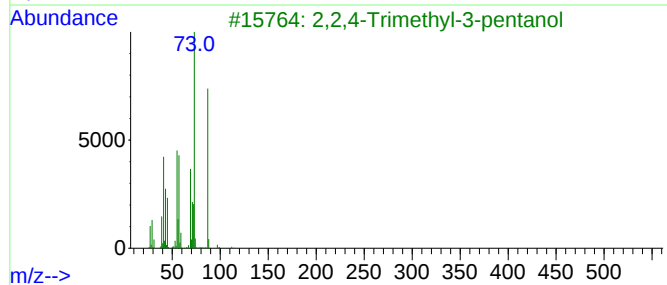
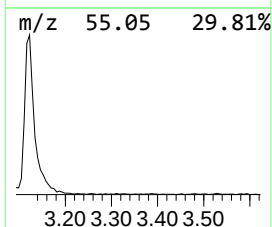
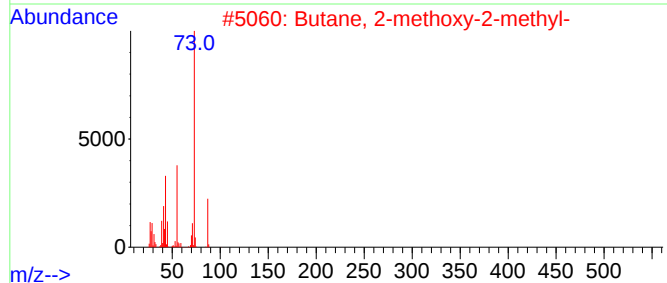
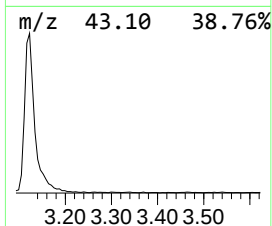
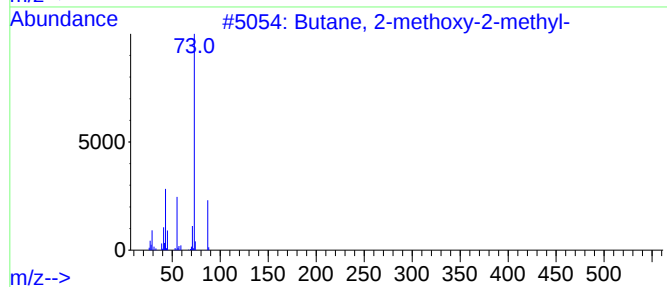
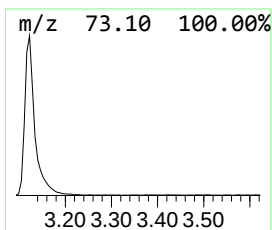
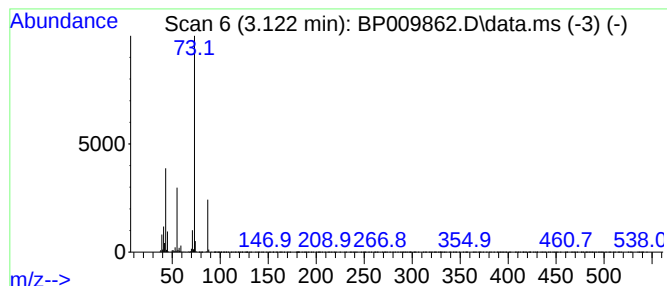
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Quant Title : SVOA CALIBRATION

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.
3.122	13.20 ng/ul	465123	1,4-Dichlorobenzene-d4	7.957

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			1,3-Dioxolane, 2-propyl-	116	C6H12O2	003390-13-4	23
5			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12



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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...	3.122	13.2	ng/u1	465123	1	7.957	704970	20.0