

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092322\
 Data File : BM036713.D
 Acq On : 24 Sep 2022 13:57
 Operator : CG/JU
 Sample : PB147750BL
 Misc :
 ALS Vial : 46 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK750

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_M\Methods\SFAM-EPA-BM091522.M
 Title : SVOA CALIBRATION

Signal : TIC: BM036713.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.116	3	5	15	rVB	225840	260934	3.86%	0.327%
2	3.387	46	51	58	rVB	95560	131685	1.95%	0.165%
3	3.816	120	124	139	rBV	869541	1365383	20.21%	1.709%
4	4.146	177	180	187	rVB	28930	40172	0.59%	0.050%
5	7.163	685	693	703	rBV	1563677	2381590	35.25%	2.981%
6	7.334	716	722	734	rVB	1306101	2013370	29.80%	2.520%
7	7.534	748	756	766	rBV	1582186	2430945	35.98%	3.043%
8	8.004	829	836	845	rBV	1463400	2288756	33.88%	2.865%
9	8.716	949	957	965	rBV	1920192	3053551	45.20%	3.823%
10	9.175	1027	1035	1043	rBV	1427253	2356414	34.88%	2.950%
11	9.898	1149	1158	1169	rBV	1067496	1767771	26.17%	2.213%
12	10.433	1242	1249	1259	rBV	1853106	3029125	44.84%	3.792%
13	10.816	1307	1314	1324	rBV	2044443	3419031	50.61%	4.280%
14	10.951	1330	1337	1353	rBV	944328	1661959	24.60%	2.081%
15	12.074	1526	1528	1538	rVB6	9193	19559	0.29%	0.024%
16	12.398	1578	1583	1591	rBV2	35466	62929	0.93%	0.079%
17	12.610	1615	1619	1623	rBV7	4872	7824	0.12%	0.010%
18	13.533	1772	1776	1783	rBV5	18138	27209	0.40%	0.034%
19	13.610	1783	1789	1791	rVV6	5205	7415	0.11%	0.009%
20	14.045	1856	1863	1878	rBV	3256700	4336095	64.19%	5.428%
21	14.327	1903	1911	1921	rBV	3701868	5196499	76.92%	6.505%
22	14.633	1955	1963	1972	rBV	3138068	4501503	66.63%	5.635%
23	14.821	1988	1995	2014	rBV	1260759	1912336	28.31%	2.394%
24	15.621	2124	2131	2143	rBV	4668973	6620919	98.01%	8.289%
25	15.733	2143	2150	2162	rVV	788089	1139943	16.87%	1.427%
26	15.857	2166	2171	2180	rVB3	40016	65889	0.98%	0.082%
27	16.292	2239	2245	2246	rBV6	5321	7310	0.11%	0.009%
28	16.398	2259	2263	2270	rVB8	11260	19307	0.29%	0.024%
29	16.856	2336	2341	2345	rVB3	12954	19192	0.28%	0.024%
30	17.051	2371	2374	2378	rBV6	9119	12967	0.19%	0.016%
31	17.156	2389	2392	2396	rVB5	6744	9454	0.14%	0.012%
32	17.368	2422	2428	2437	rBV	3441550	4664596	69.05%	5.839%
33	17.468	2438	2445	2465	rVV	5068336	6755571	100.00%	8.457%
34	17.933	2520	2524	2527	rVB4	12398	17505	0.26%	0.022%
35	18.039	2538	2542	2545	rBV5	11406	13273	0.20%	0.017%
36	19.315	2756	2759	2764	rVB	47489	57791	0.86%	0.072%

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 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 >
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37	19.386	2767	2771	2779	rBV2	49867	79234	1.17%	0.099%
38	19.750	2827	2833	2842	rBV	4968085	6520190	96.52%	8.162%
39	20.744	2999	3002	3005	rBV2	148975	158704	2.35%	0.199%
40	21.533	3132	3136	3143	rBV	2715294	3431974	50.80%	4.296%
41	23.809	3517	3523	3532	rBV2	2396065	4838783	71.63%	6.058%
42	23.968	3544	3550	3560	rVB2	1584465	3175994	47.01%	3.976%

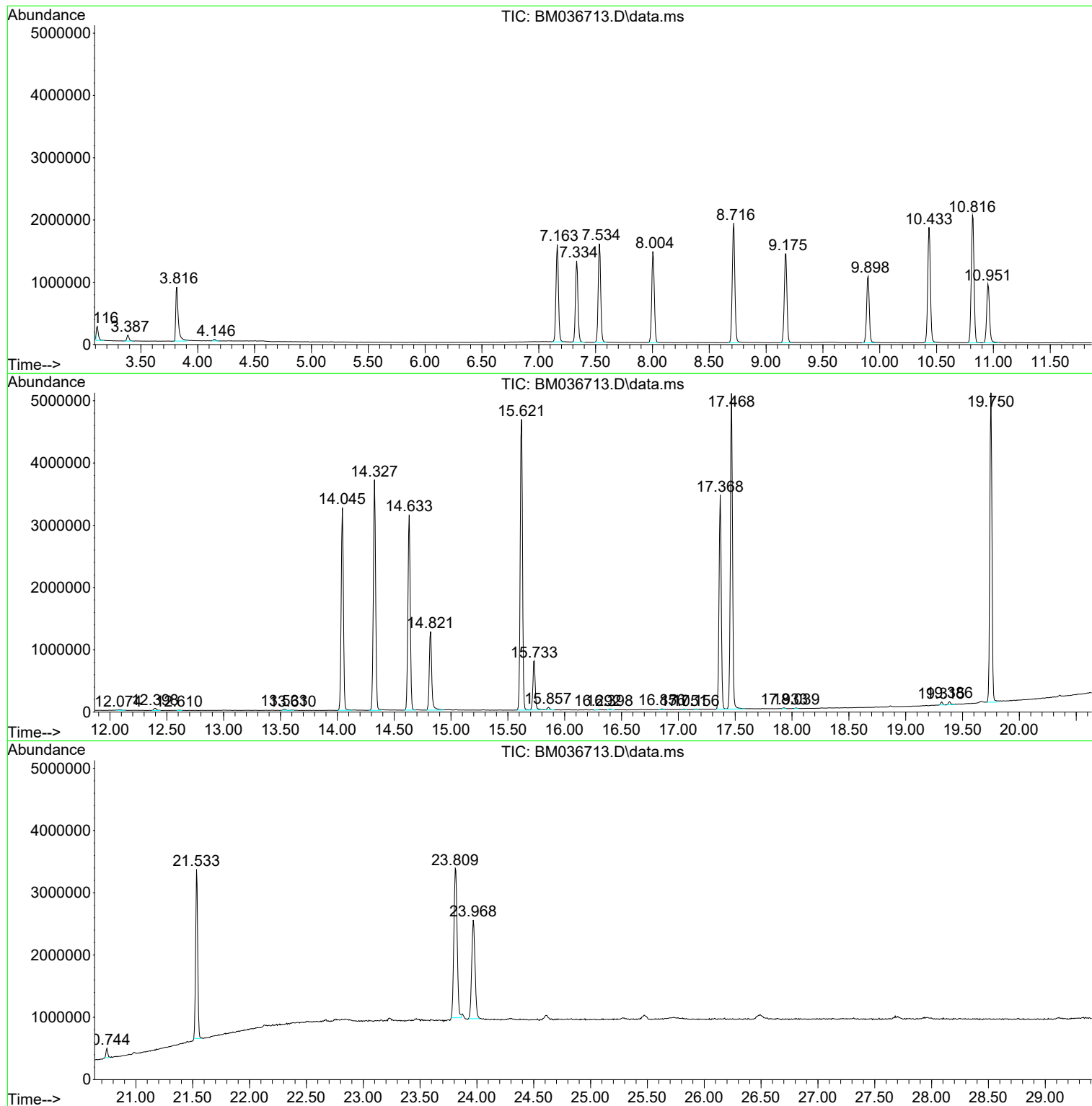
Sum of corrected areas: 79880651

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
Quant Title : SVOA CALIBRATION

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



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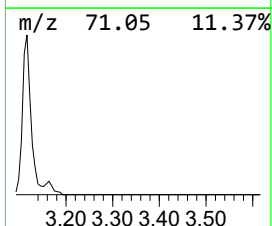
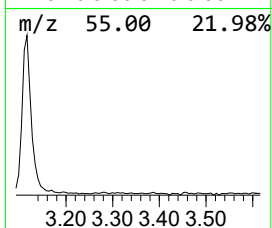
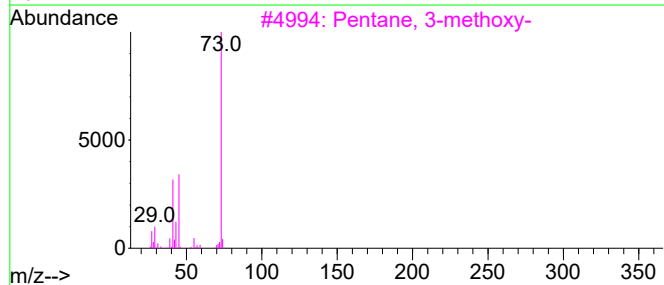
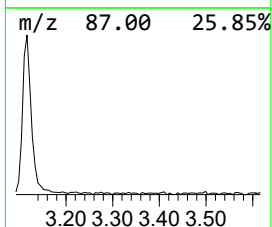
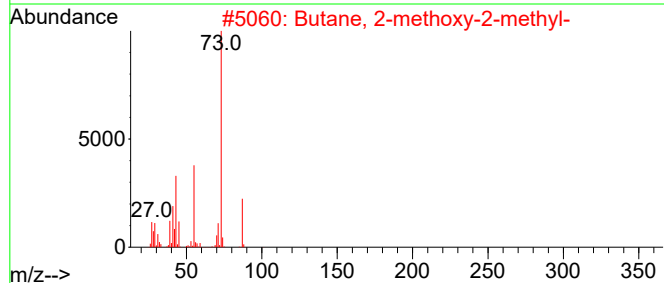
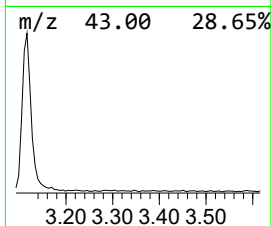
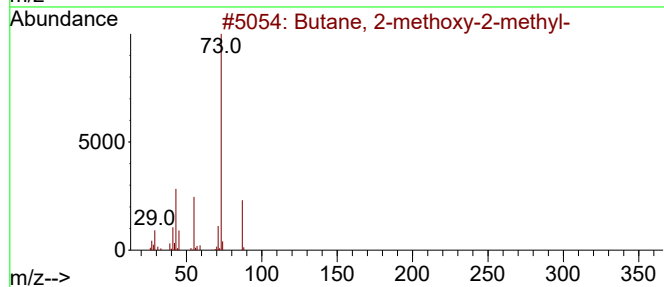
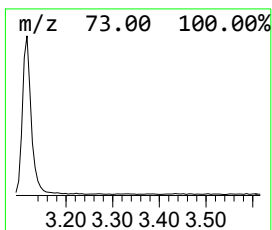
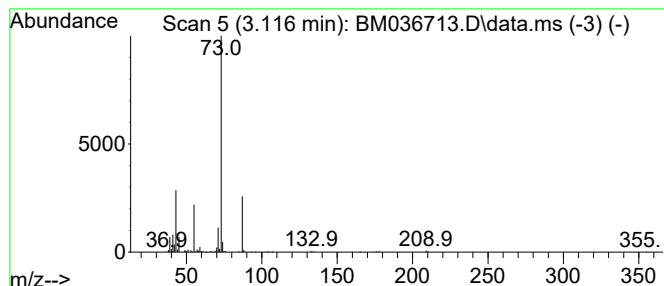
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Quant Title : SVOA CALIBRATION

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.
3.116	2.28 ng/ul	260934	1,4-Dichlorobenzene-d4	8.004

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	28		
3	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17		
4	Silane, tetramethyl-	88	C4H12Si	000075-76-3	12		
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...	3.116	2.3	ng/ul	260934	1	8.004	2288760	20.0