

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041923\
 Data File : BP014724.D
 Acq On : 19 Apr 2023 11:41
 Operator : CG/JU
 Sample : 02336-08
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 COMG7

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-BP041823.M
 Title : SVOA CALIBRATION

Signal : TIC: BP014724.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.217	3	5	20	rVB	1096767	1173066	20.88%	2.462%
2	3.487	47	51	57	rBV	49913	60550	1.08%	0.127%
3	3.922	124	125	141	rBV	116585	173969	3.10%	0.365%
4	4.164	162	166	170	rVB2	7913	10404	0.19%	0.022%
5	4.264	179	183	188	rBV	16772	24698	0.44%	0.052%
6	5.205	339	343	349	rVB4	4928	7164	0.13%	0.015%
7	7.287	691	697	707	rVB	199025	305967	5.45%	0.642%
8	7.469	721	728	739	rBV	710991	1082819	19.28%	2.272%
9	7.675	756	763	776	rBV	652631	1013720	18.05%	2.127%
10	7.769	776	779	784	rVB6	4509	5683	0.10%	0.012%
11	8.152	836	844	852	rBV	686085	1105126	19.67%	2.319%
12	8.846	955	962	976	rBV	527998	842438	15.00%	1.768%
13	9.322	1033	1043	1051	rBV	690800	1104025	19.65%	2.317%
14	9.740	1109	1114	1119	rVB7	3998	7726	0.14%	0.016%
15	10.051	1159	1167	1178	rBV	420597	669346	11.92%	1.405%
16	10.587	1248	1258	1268	rBV	859384	1406867	25.05%	2.953%
17	10.981	1316	1325	1333	rBV	922651	1540095	27.42%	3.232%
18	11.110	1336	1347	1359	rBV	594380	986230	17.56%	2.070%
19	12.493	1577	1582	1587	rVB8	4023	7419	0.13%	0.016%
20	12.551	1587	1592	1599	rVB3	17669	28346	0.50%	0.059%
21	13.675	1779	1783	1787	rBV6	5682	8662	0.15%	0.018%
22	14.187	1862	1870	1882	rBV	1821058	2518858	44.84%	5.286%
23	14.310	1889	1891	1897	rVB7	4475	6938	0.12%	0.015%
24	14.481	1912	1920	1929	rBV	2119473	3021339	53.79%	6.341%
25	14.787	1963	1972	1980	rBV	1382172	2057399	36.63%	4.318%
26	14.845	1980	1982	1988	rVB7	3998	6459	0.11%	0.014%
27	14.945	1993	1999	2007	rBV	121065	183219	3.26%	0.385%
28	15.187	2036	2040	2045	rVB8	4858	8160	0.15%	0.017%
29	15.775	2132	2140	2147	rBV	2736484	3847596	68.50%	8.075%
30	15.875	2151	2157	2168	rBV	375261	529708	9.43%	1.112%
31	16.010	2176	2180	2187	rVB3	11070	18750	0.33%	0.039%
32	16.134	2196	2201	2207	rBV	125825	169831	3.02%	0.356%
33	17.533	2432	2439	2446	rBV	1746378	2461885	43.83%	5.167%
34	17.633	2449	2456	2471	rVV	3192169	4374925	77.89%	9.181%
35	18.392	2581	2585	2593	rBV5	27058	49073	0.87%	0.103%
36	19.427	2756	2761	2766	rBV2	38535	55032	0.98%	0.115%

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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-BP041823.M
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37	19.492	2768	2772	2776	rVV	36026	48227	0.86%	0.101%
38	19.622	2791	2794	2796	rBV4	13200	15195	0.27%	0.032%
39	19.851	2826	2833	2842	rBV	3812126	5253863	93.53%	11.026%
40	21.286	3074	3077	3082	rBV2	46073	54109	0.96%	0.114%
41	21.604	3126	3131	3142	rBV	2041544	2768938	49.29%	5.811%
42	24.021	3534	3542	3559	rVV	2858915	5617128	100.00%	11.788%
43	24.186	3562	3570	3583	rVB2	1418133	3018670	53.74%	6.335%

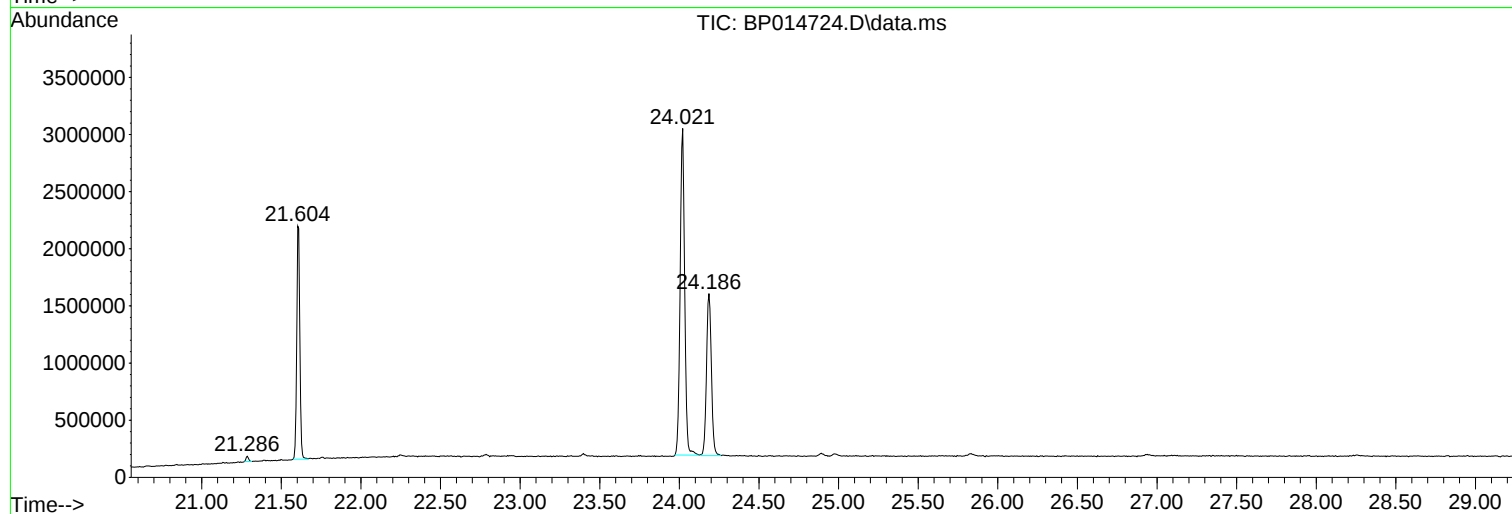
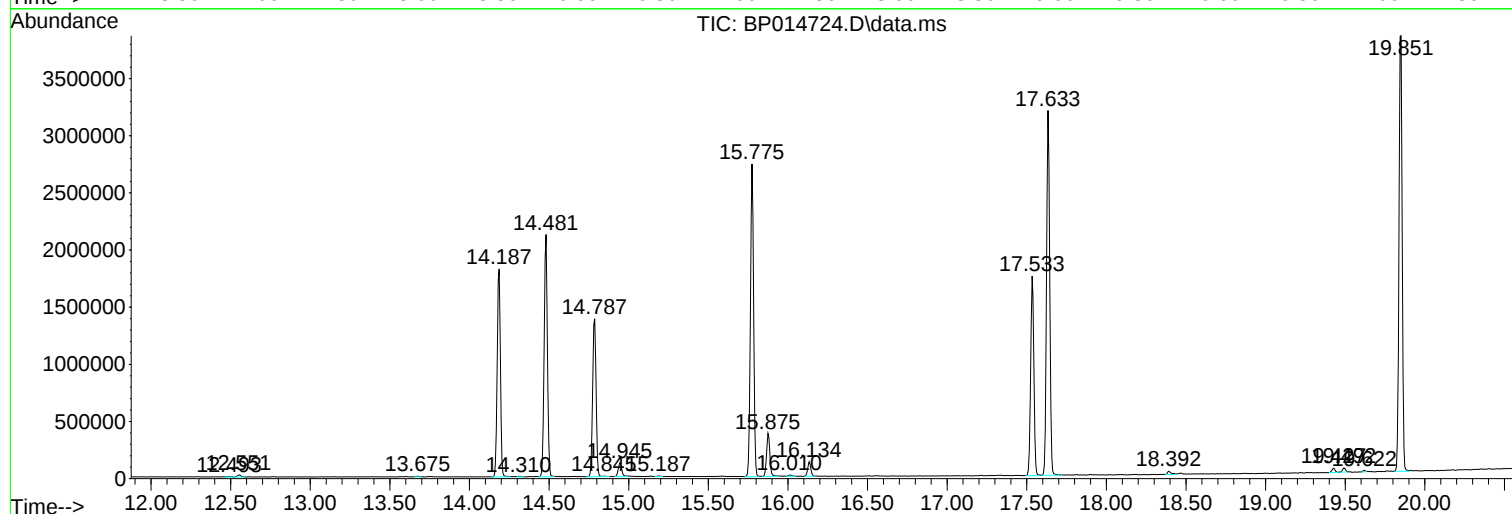
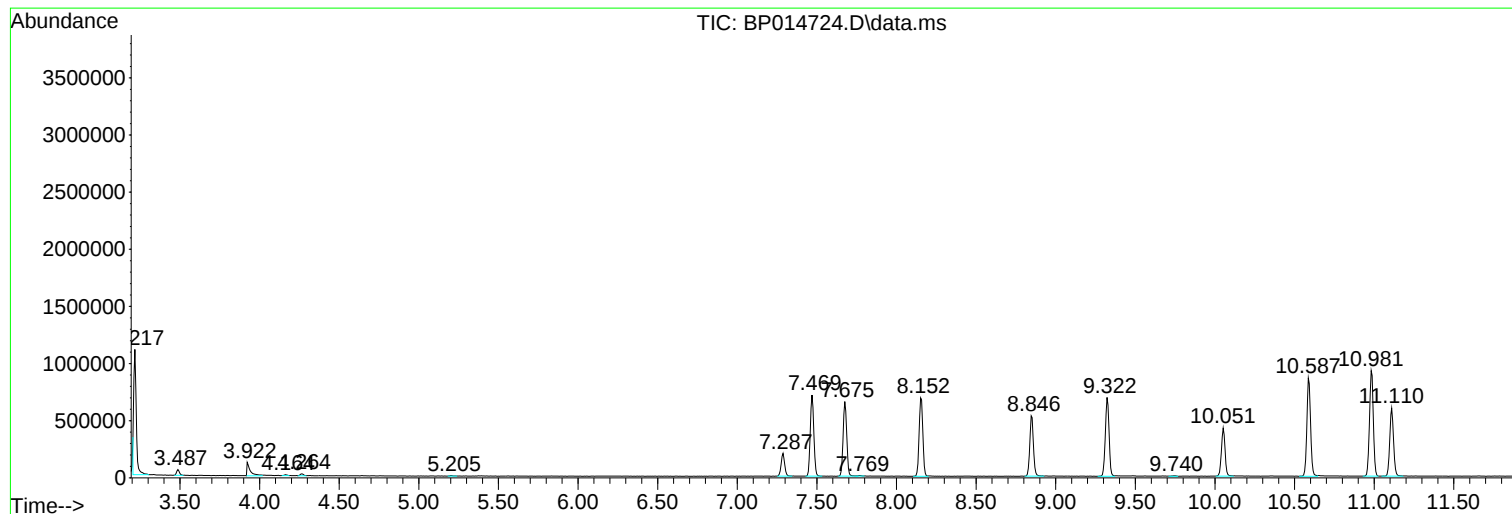
Sum of corrected areas: 47649622

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP041823.M
Quant Title : SVOA CALIBRATION

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



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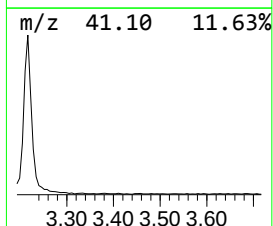
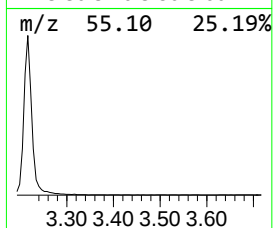
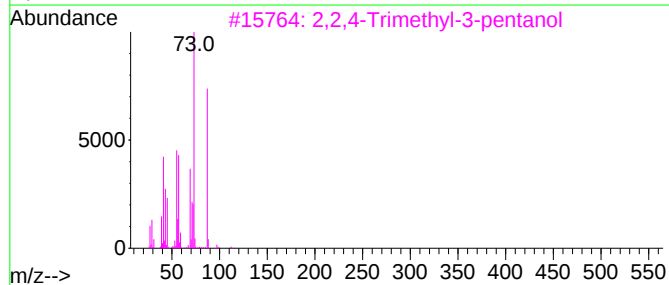
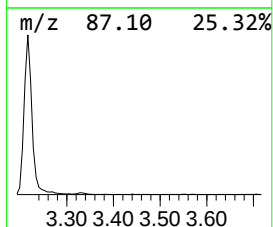
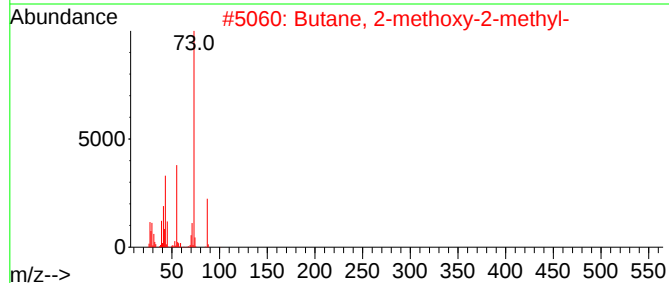
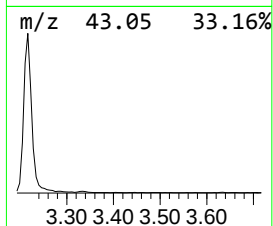
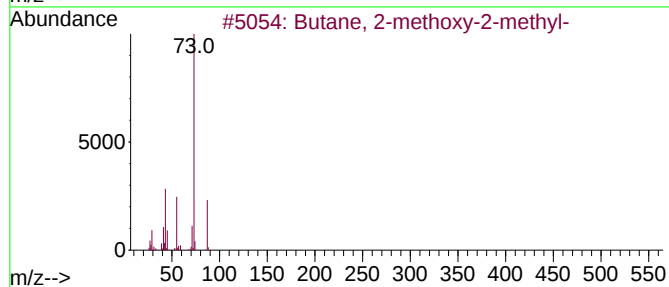
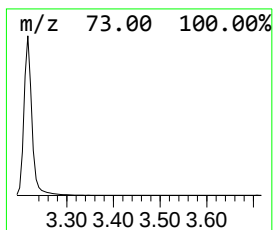
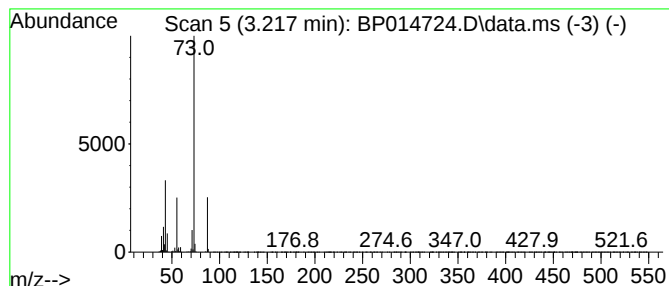
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TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.217	21.23 ng/ul	1173070	1,4-Dichlorobenzene-d4	8.152

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	74		
3	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39		
4	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23		
5	Silane, tetramethyl-	88	C4H12Si	000075-76-3	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...	3.217	21.2	ng/u1	1173070	1	8.152	1105130	20.0