

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042023\
 Data File : BP014781.D
 Acq On : 21 Apr 2023 09:35
 Operator : CG/JU
 Sample : 02373-01
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 CB9D1

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: BP014781.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	21	rVB	1024631	1127537	20.05%	1.880%
2	3.487	47	51	56	rVB	59990	74358	1.32%	0.124%
3	3.923	123	125	140	rBV	272504	404598	7.19%	0.675%
4	4.034	140	144	152	rVB10	8837	18665	0.33%	0.031%
5	4.164	163	166	171	rVB4	6754	11172	0.20%	0.019%
6	4.270	179	184	189	rVB	23812	35674	0.63%	0.059%
7	4.487	217	221	225	rBV4	14410	18266	0.32%	0.030%
8	4.911	289	293	298	rVB7	7010	9810	0.17%	0.016%
9	5.205	335	343	348	rBV8	5233	10215	0.18%	0.017%
10	5.393	370	375	380	rBV	20661	31481	0.56%	0.052%
11	6.522	562	567	570	rBV7	4859	7367	0.13%	0.012%
12	7.287	691	697	706	rVB	229154	358375	6.37%	0.598%
13	7.469	719	728	739	rBV	868707	1335693	23.75%	2.227%
14	7.675	756	763	773	rBV	812502	1225242	21.79%	2.043%
15	8.152	837	844	851	rBV	1061156	1651661	29.37%	2.754%
16	8.852	955	963	975	rBV	607499	984110	17.50%	1.641%
17	9.322	1035	1043	1053	rBV	908969	1464785	26.05%	2.442%
18	9.411	1053	1058	1060	rVB6	4166	6566	0.12%	0.011%
19	9.734	1108	1113	1118	rVB5	6591	11388	0.20%	0.019%
20	10.052	1159	1167	1179	rBV	635878	1008943	17.94%	1.682%
21	10.199	1188	1192	1197	rVB5	7744	11671	0.21%	0.019%
22	10.587	1250	1258	1271	rBV	1131150	1828602	32.52%	3.049%
23	10.981	1316	1325	1333	rBV	1523974	2574607	45.78%	4.293%
24	11.110	1337	1347	1361	rBV	587339	992548	17.65%	1.655%
25	11.634	1431	1436	1443	rBV6	4970	10629	0.19%	0.018%
26	11.787	1457	1462	1467	rBV3	21329	34080	0.61%	0.057%
27	12.552	1587	1592	1599	rVB2	24279	37640	0.67%	0.063%
28	13.599	1767	1770	1776	rVB8	5965	8549	0.15%	0.014%
29	13.669	1776	1782	1790	rBV6	7439	12672	0.23%	0.021%
30	14.181	1862	1869	1880	rBV	2438424	3371505	59.95%	5.622%
31	14.304	1886	1890	1896	rVB9	6655	10758	0.19%	0.018%
32	14.475	1913	1919	1929	rBV	2549871	3774198	67.11%	6.293%
33	14.663	1949	1951	1956	rVB6	5640	6539	0.12%	0.011%
34	14.781	1963	1971	1979	rBV	2430193	3520615	62.60%	5.870%
35	14.846	1979	1982	1989	rVB9	6409	10663	0.19%	0.018%
36	14.946	1993	1999	2006	rBV	170044	246359	4.38%	0.411%

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Peak Location: TOP

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP041823.M
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37	15.110	2023	2027	2030	rBV6	6235	8519	0.15%	0.014%
38	15.175	2034	2038	2043	rVB8	7708	12419	0.22%	0.021%
39	15.775	2132	2140	2151	rBV	3393906	5047370	89.75%	8.416%
40	15.875	2151	2157	2171	rVB	763781	993046	17.66%	1.656%
41	16.016	2177	2181	2186	rVB2	14974	21992	0.39%	0.037%
42	16.134	2195	2201	2207	rBV	195510	264101	4.70%	0.440%
43	16.557	2269	2273	2278	rVB7	4286	6639	0.12%	0.011%
44	16.910	2329	2333	2336	rBV6	5330	6736	0.12%	0.011%
45	17.204	2380	2383	2391	rVB9	5869	10947	0.19%	0.018%
46	17.275	2391	2395	2399	rBV7	6495	13167	0.23%	0.022%
47	17.534	2432	2439	2449	rBV	2796545	3854083	68.53%	6.426%
48	17.634	2449	2456	2466	rVV	3699600	5170307	91.94%	8.621%
49	17.734	2471	2473	2479	rVB7	7579	10950	0.19%	0.018%
50	17.881	2494	2498	2505	rBV8	14828	29082	0.52%	0.048%
51	18.257	2560	2562	2568	rVB7	8390	10977	0.20%	0.018%
52	18.381	2579	2583	2593	rBV	148409	216272	3.85%	0.361%
53	18.716	2637	2640	2645	rVB4	17023	22903	0.41%	0.038%
54	19.422	2756	2760	2764	rBV3	47450	61634	1.10%	0.103%
55	19.486	2765	2771	2775	rBV	49679	97187	1.73%	0.162%
56	19.604	2787	2791	2796	rBV4	87338	109944	1.96%	0.183%
57	19.845	2826	2832	2842	rBV	4455256	5623642	100.00%	9.377%
58	21.275	3072	3075	3083	rBV2	178390	244286	4.34%	0.407%
59	21.604	3125	3131	3137	rBV	2447560	3277977	58.29%	5.466%
60	22.922	3352	3355	3361	rVB	115265	164465	2.92%	0.274%
61	24.010	3532	3540	3556	rVB2	2433596	5103766	90.76%	8.510%
62	24.174	3561	3568	3579	rVB2	1594255	3343674	59.46%	5.575%

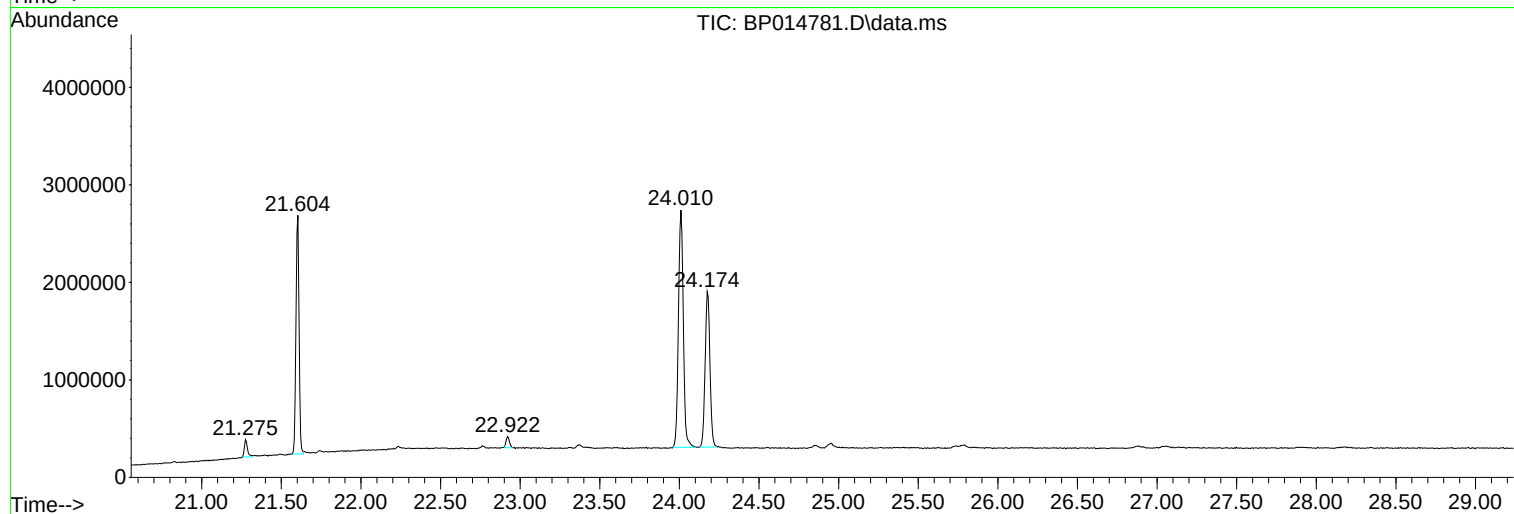
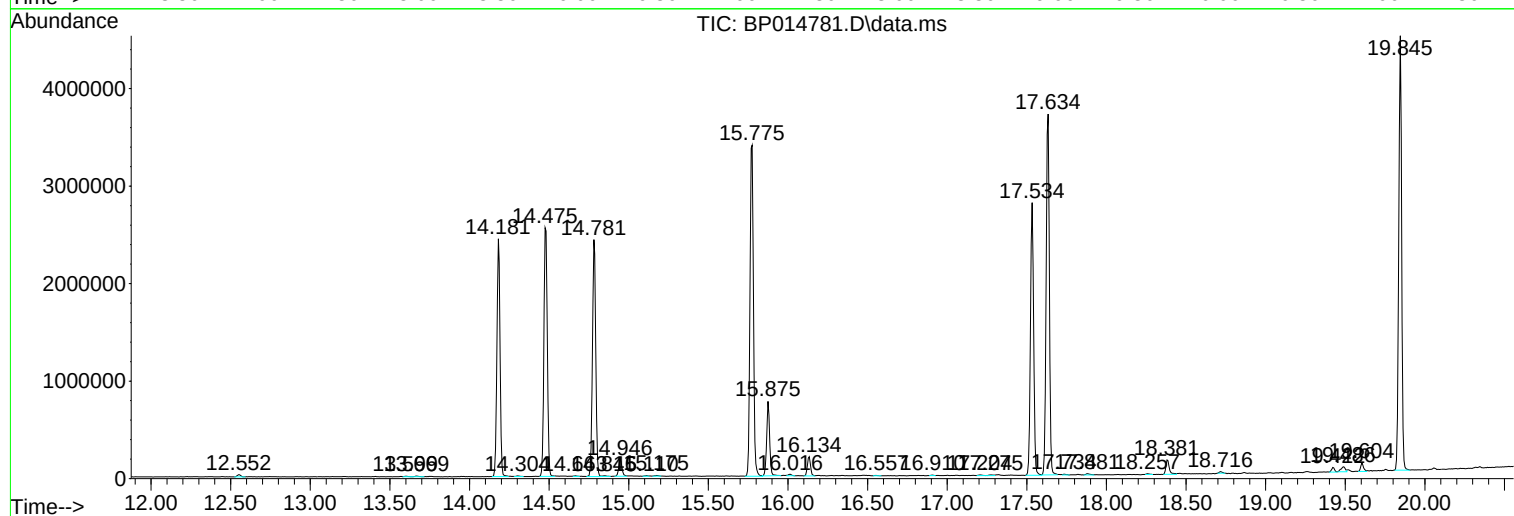
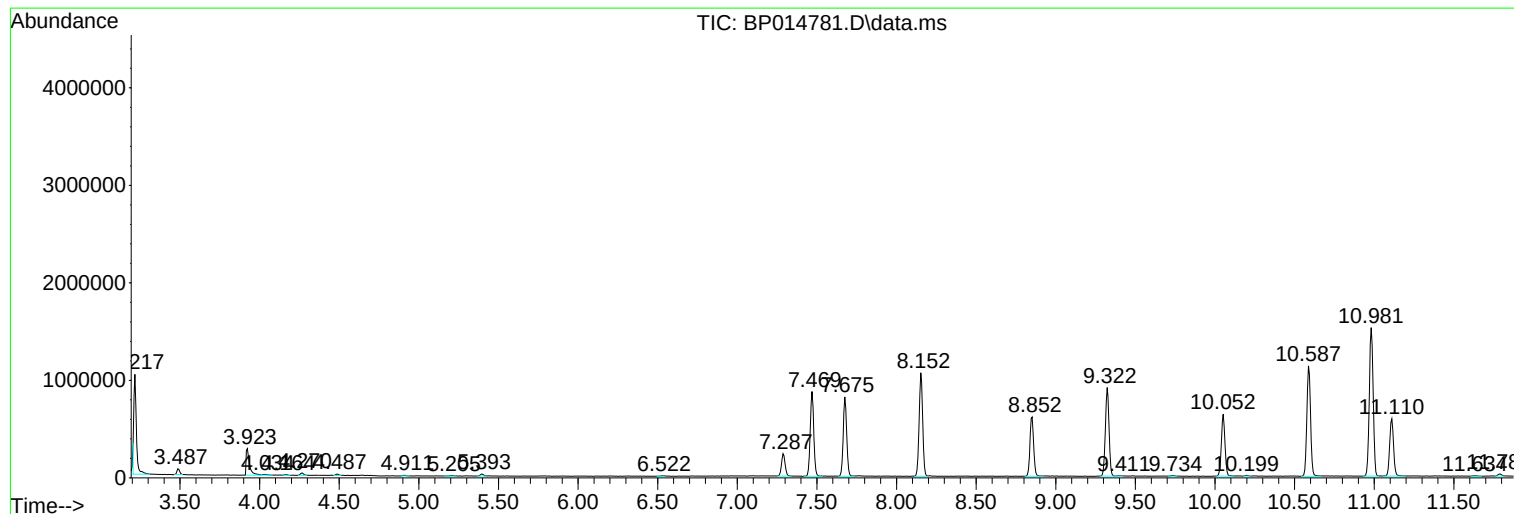
Sum of corrected areas: 59973626

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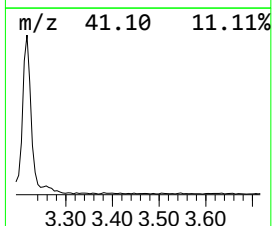
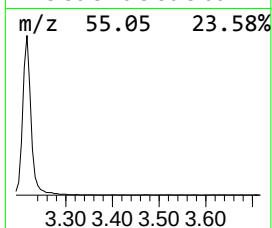
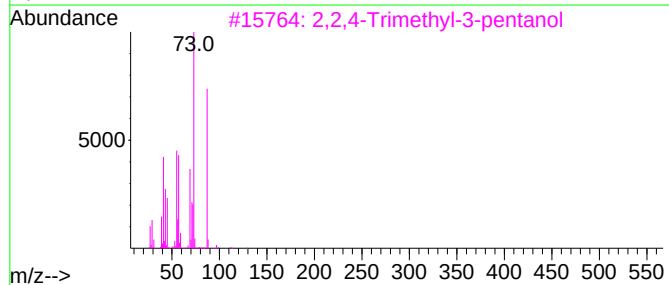
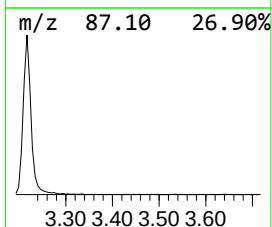
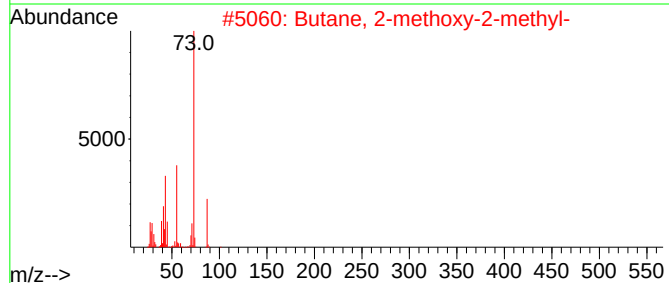
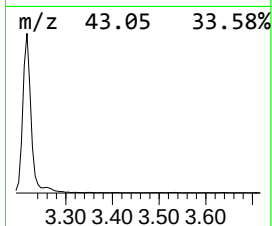
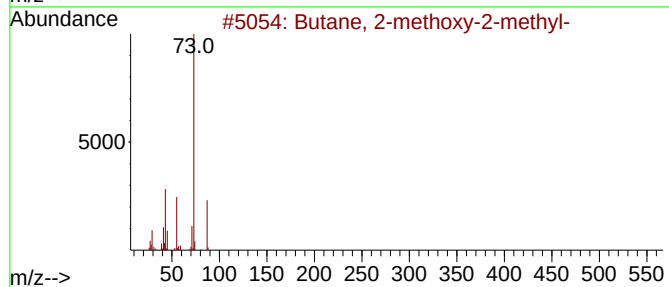
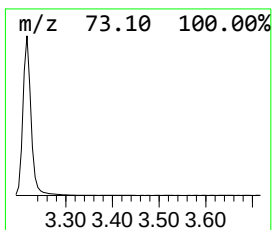
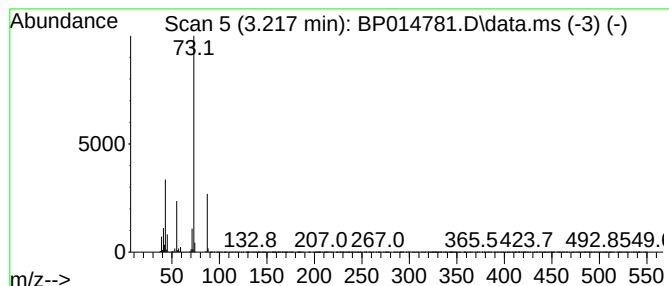
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP041823.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.217	13.65 ng/ul	1127540	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	83		
2	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	64		
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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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	13.7	ng/ul	1127540	1	8.152	1651660	20.0