

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP083024\
 Data File : BP021773.D
 Acq On : 31 Aug 2024 06:06
 Operator : RC/JU
 Sample : P3733-01
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A4495

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-BP082324.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP021773.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.017	3	5	19	rVB	3745751	4886472	93.63%	7.566%
2	3.276	45	49	55	rBV	67619	87436	1.68%	0.135%
3	3.546	91	95	103	rBV	60894	86435	1.66%	0.134%
4	3.687	115	119	130	rBV	835367	1204551	23.08%	1.865%
5	4.017	171	175	179	rVB	21015	28067	0.54%	0.043%
6	4.334	224	229	234	rBV	53630	69214	1.33%	0.107%
7	4.811	305	310	315	rVB8	5673	9967	0.19%	0.015%
8	4.928	324	330	335	rBV2	67736	100661	1.93%	0.156%
9	6.058	516	522	526	rVB3	11403	16920	0.32%	0.026%
10	6.787	641	646	659	rVB4	9970	22908	0.44%	0.035%
11	6.952	667	674	684	rBV	987994	1576955	30.22%	2.442%
12	7.111	693	701	712	rVB	802117	1293472	24.78%	2.003%
13	7.322	729	737	745	rBV	948908	1546707	29.64%	2.395%
14	7.393	745	749	757	rVB2	17530	35635	0.68%	0.055%
15	7.781	808	815	823	rBV	1182081	1990988	38.15%	3.083%
16	8.158	874	879	880	rBV5	4162	5705	0.11%	0.009%
17	8.487	927	935	950	rBV	1089956	1873583	35.90%	2.901%
18	8.928	1001	1010	1026	rBV	869988	1413441	27.08%	2.188%
19	9.052	1028	1031	1035	rVV7	6153	9507	0.18%	0.015%
20	9.099	1035	1039	1045	rVB8	5800	9652	0.18%	0.015%
21	9.493	1101	1106	1112	rVB3	13875	21413	0.41%	0.033%
22	9.652	1125	1133	1142	rBV	594941	1043614	20.00%	1.616%
23	10.187	1213	1224	1233	rBV	841076	1569746	30.08%	2.430%
24	10.563	1280	1288	1297	rBV	1507723	2532218	48.52%	3.921%
25	10.699	1304	1311	1331	rVB	903665	1653312	31.68%	2.560%
26	11.116	1374	1382	1386	rBV10	6521	16766	0.32%	0.026%
27	11.310	1409	1415	1431	rBV	52229	143107	2.74%	0.222%
28	12.099	1544	1549	1552	rBV5	4524	7369	0.14%	0.011%
29	12.163	1554	1560	1563	rBV5	13298	25804	0.49%	0.040%
30	12.775	1660	1664	1669	rBV8	6039	10739	0.21%	0.017%
31	13.028	1700	1707	1713	rVB9	7605	14161	0.27%	0.022%
32	13.322	1755	1757	1763	rVB7	4741	7876	0.15%	0.012%
33	13.387	1763	1768	1774	rBV10	4343	8459	0.16%	0.013%
34	13.610	1800	1806	1808	rBV7	4078	6018	0.12%	0.009%
35	13.828	1835	1843	1857	rBV	1549621	2225832	42.65%	3.446%
36	14.110	1881	1891	1910	rBV	1740501	2733865	52.38%	4.233%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
 Title : SVOA CALIBRATION

37	14.428	1935	1945	1958	rBV	1986147	3120963	59.80%	4.832%
38	14.628	1973	1979	2003	rBV	525504	1178769	22.59%	1.825%
39	14.851	2013	2017	2022	rVB8	9360	16265	0.31%	0.025%
40	15.251	2080	2085	2088	rBV4	7417	13386	0.26%	0.021%
41	15.434	2108	2116	2130	rBV	2247934	3503448	67.13%	5.424%
42	15.551	2130	2136	2145	rVV	476819	720250	13.80%	1.115%
43	15.675	2154	2157	2164	rVB7	10867	16650	0.32%	0.026%
44	16.016	2210	2215	2219	rBV7	5973	10636	0.20%	0.016%
45	16.181	2239	2243	2249	rBV9	6732	10497	0.20%	0.016%
46	17.004	2378	2383	2391	rBV8	11965	31667	0.61%	0.049%
47	17.228	2413	2421	2429	rBV	2265130	3713115	71.15%	5.749%
48	17.334	2432	2439	2451	rBV2	2408786	3657787	70.09%	5.663%
49	17.422	2451	2454	2460	rVB7	22432	36885	0.71%	0.057%
50	17.940	2539	2542	2549	rBV6	13465	19949	0.38%	0.031%
51	18.169	2576	2581	2590	rBV	154525	271317	5.20%	0.420%
52	19.251	2763	2765	2771	rVB7	34333	39601	0.76%	0.061%
53	19.328	2775	2778	2782	rBV3	33714	43904	0.84%	0.068%
54	19.504	2804	2808	2813	rBV6	52963	79914	1.53%	0.124%
55	19.698	2835	2841	2851	rVB	3169426	4663577	89.36%	7.220%
56	21.363	3119	3124	3130	rBV2	234053	407943	7.82%	0.632%
57	21.675	3172	3177	3184	rBV2	2939865	4575747	87.67%	7.084%
58	24.839	3707	3715	3725	rVB	1839854	4948361	94.81%	7.661%
59	25.069	3746	3754	3768	rVB	1796639	5218988	100.00%	8.080%

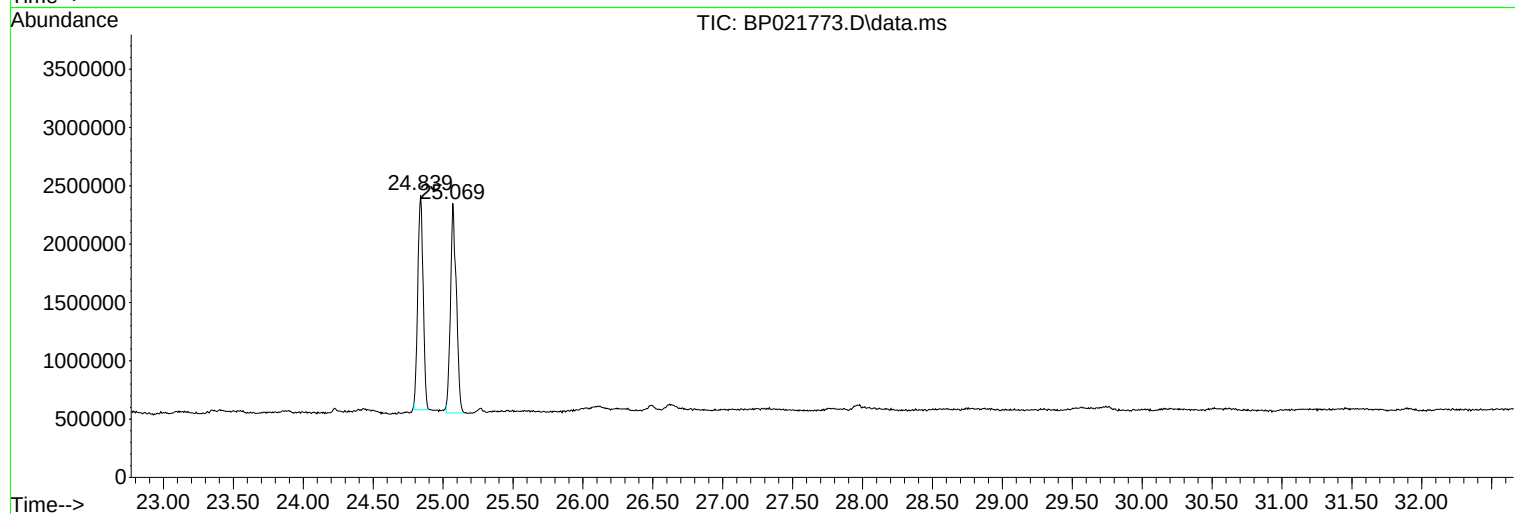
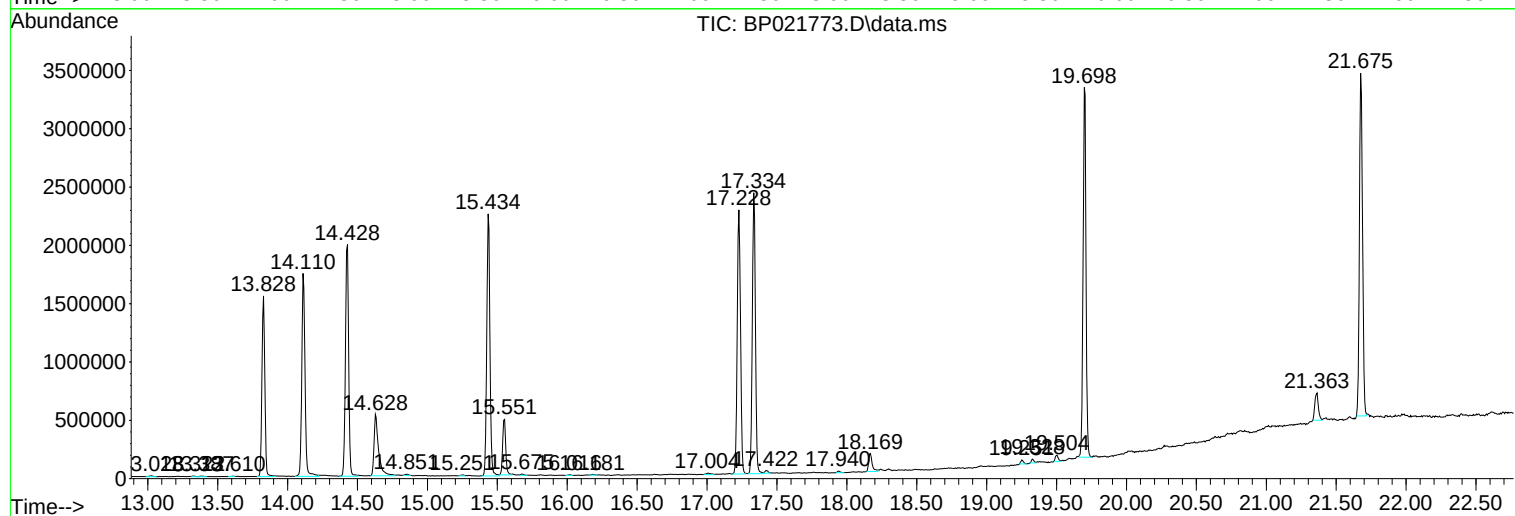
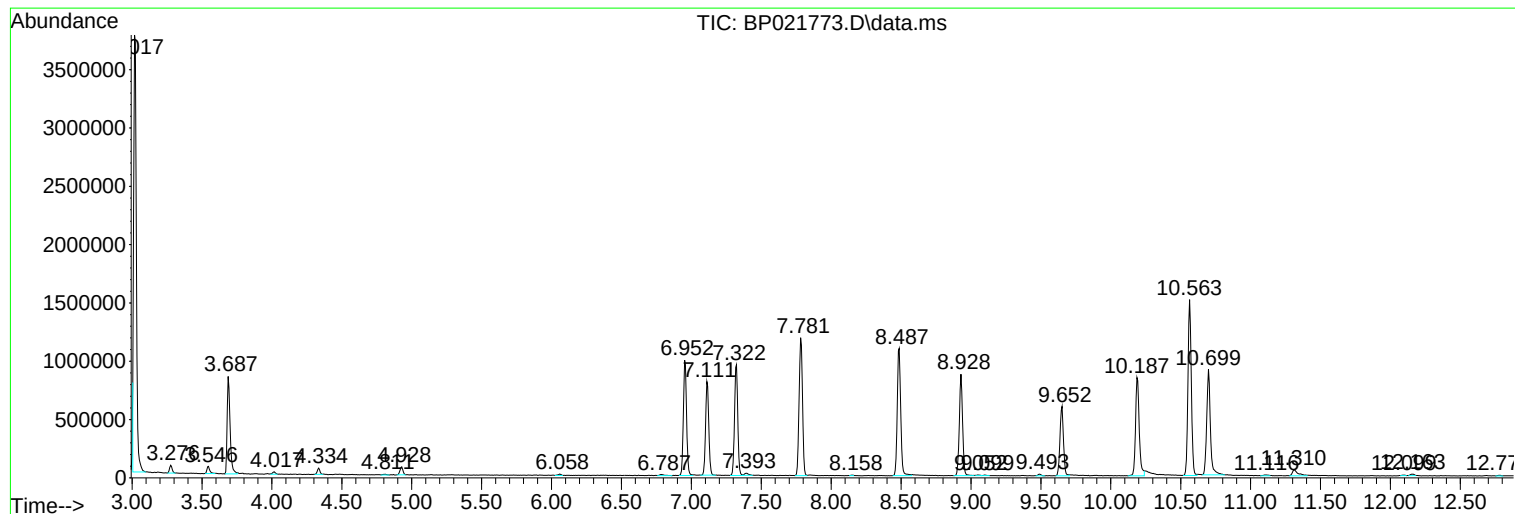
Sum of corrected areas: 64588194

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

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



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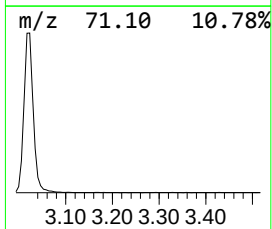
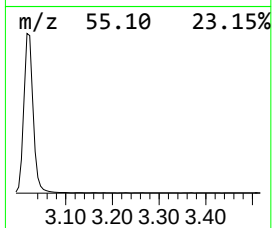
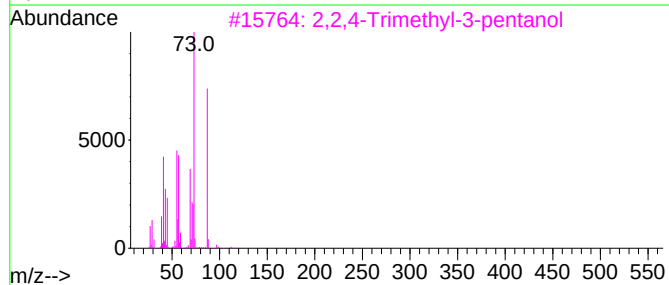
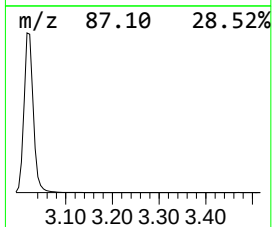
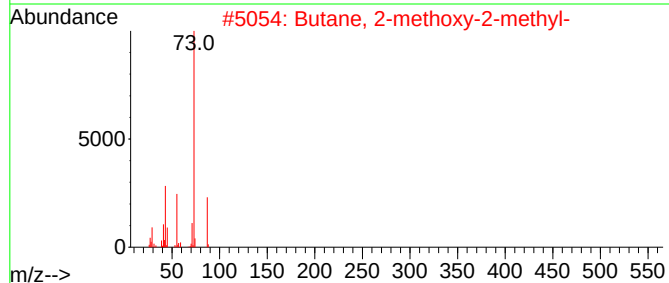
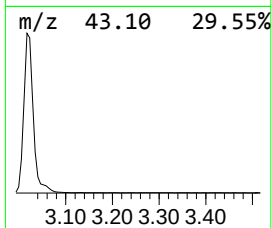
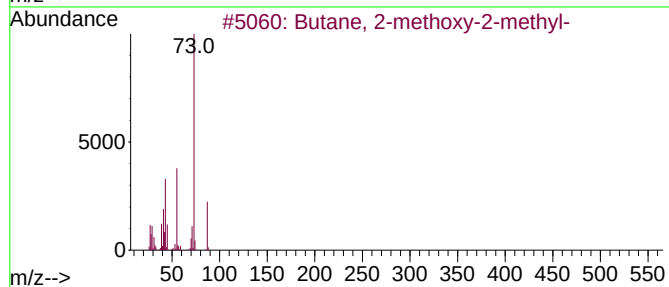
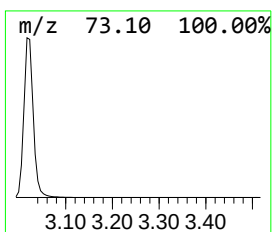
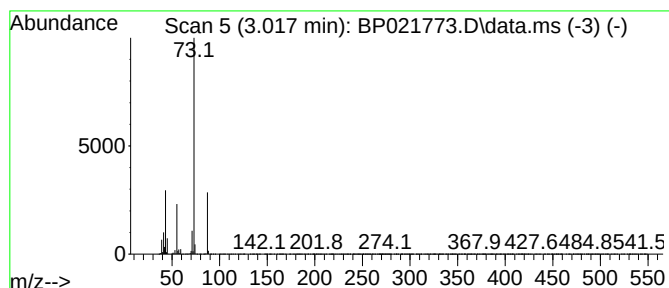
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

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.
3.017	49.09 ng/ul	4886470	1,4-Dichlorobenzene-d4	7.781

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	72		
3	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	43		
4	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23		
5	Borane, dimethoxy-	74	C2H7BO2	004542-61-4	9		



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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-metho...	3.017	49.1	ng/ul	4886470	1	7.781	1990990	20.0