

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP083024\
 Data File : BP021770.D
 Acq On : 31 Aug 2024 03:55
 Operator : RC/JU
 Sample : PB163036BL
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK036

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: BP021770.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.022	3	6	17	rVB	162940	217955	2.99%	0.287%
2	3.116	18	22	28	rVB	22346	35260	0.48%	0.046%
3	3.275	45	49	60	rVB	107434	150053	2.06%	0.198%
4	3.687	115	119	136	rBV	1228623	1792570	24.63%	2.364%
5	4.016	170	175	183	rVB	28031	40870	0.56%	0.054%
6	4.475	251	253	261	rVB9	4127	8124	0.11%	0.011%
7	6.957	668	675	685	rBV	1377613	2222211	30.53%	2.930%
8	7.116	693	702	715	rVB	1131912	1816002	24.95%	2.395%
9	7.322	729	737	746	rBV	1372895	2134842	29.33%	2.815%
10	7.781	808	815	823	rVB	1212270	1966533	27.02%	2.593%
11	8.487	927	935	948	rBV	1479504	2660691	36.56%	3.508%
12	8.928	1002	1010	1022	rBV	1213572	2017445	27.72%	2.660%
13	9.651	1124	1133	1143	rBV	899594	1537722	21.13%	2.028%
14	10.192	1217	1225	1233	rBV	1272563	2283923	31.38%	3.012%
15	10.569	1281	1289	1301	rBV	1449362	2530853	34.77%	3.337%
16	10.698	1303	1311	1338	rVB	1177454	2308378	31.72%	3.044%
17	11.369	1421	1425	1430	rBV7	6047	10697	0.15%	0.014%
18	12.163	1553	1560	1568	rBV	20817	40739	0.56%	0.054%
19	13.822	1836	1842	1852	rBV	2362997	3401392	46.74%	4.485%
20	14.110	1883	1891	1913	rBV	2389675	4027500	55.34%	5.311%
21	14.422	1936	1944	1955	rBV	1934707	3168743	43.54%	4.178%
22	14.627	1972	1979	1999	rBV	922255	1750061	24.05%	2.308%
23	14.845	2012	2016	2021	rVB8	7478	14744	0.20%	0.019%
24	15.433	2108	2116	2130	rBV	3204673	5066111	69.61%	6.680%
25	15.551	2130	2136	2152	rVV	979515	1433240	19.69%	1.890%
26	15.680	2154	2158	2162	rVB5	13502	22717	0.31%	0.030%
27	15.951	2200	2204	2210	rBV9	4800	7949	0.11%	0.010%
28	16.016	2210	2215	2218	rBV7	3496	7548	0.10%	0.010%
29	16.227	2247	2251	2257	rBV7	10682	22622	0.31%	0.030%
30	16.898	2359	2365	2371	rBV7	7489	13760	0.19%	0.018%
31	17.004	2380	2383	2385	rBV4	6719	8314	0.11%	0.011%
32	17.221	2413	2420	2431	rBV	2502809	3858349	53.01%	5.088%
33	17.327	2431	2438	2452	rBV2	3383484	5533593	76.03%	7.297%
34	18.157	2577	2579	2585	rVB7	9431	11930	0.16%	0.016%
35	19.251	2761	2765	2768	rBV2	45858	61057	0.84%	0.081%
36	19.327	2774	2778	2781	rVB	39285	46742	0.64%	0.062%

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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-BP082324.MA.M
Title : SVOA CALIBRATION

37	19.698	2835	2841	2853	rBV2	4504328	6862525	94.29%	9.049%
38	21.686	3173	3179	3186	rBV	2954136	4595470	63.14%	6.060%
39	24.862	3709	3719	3733	rVB2	2675951	7277888	100.00%	9.597%
40	25.098	3750	3759	3775	rVB	1537636	4870892	66.93%	6.423%

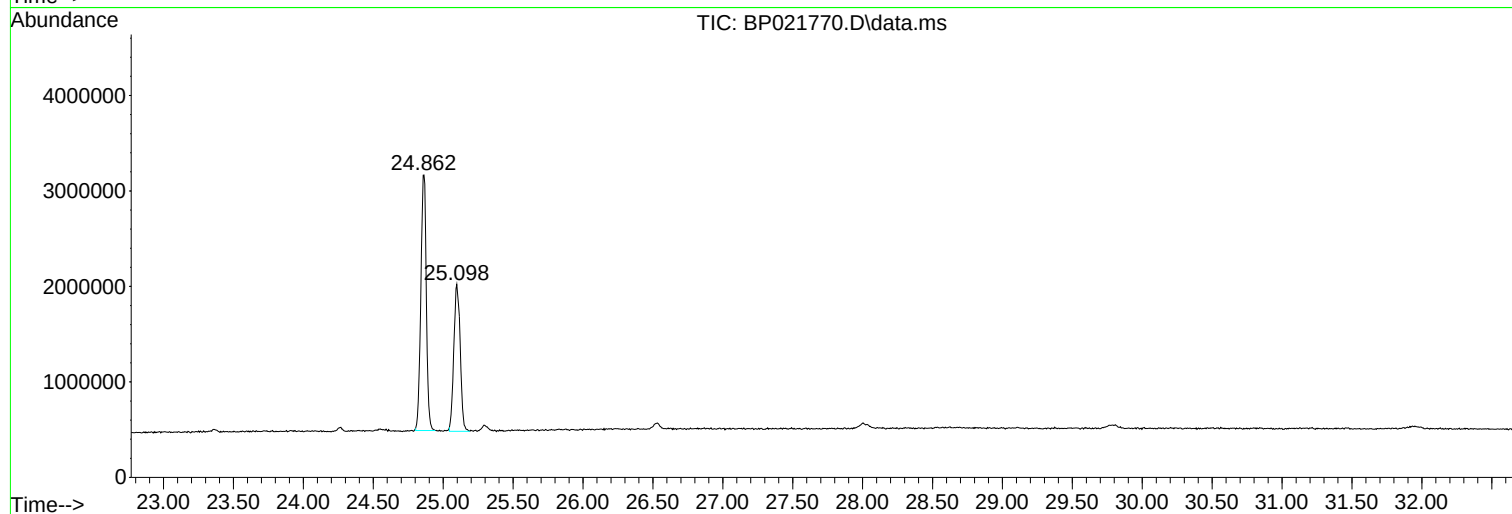
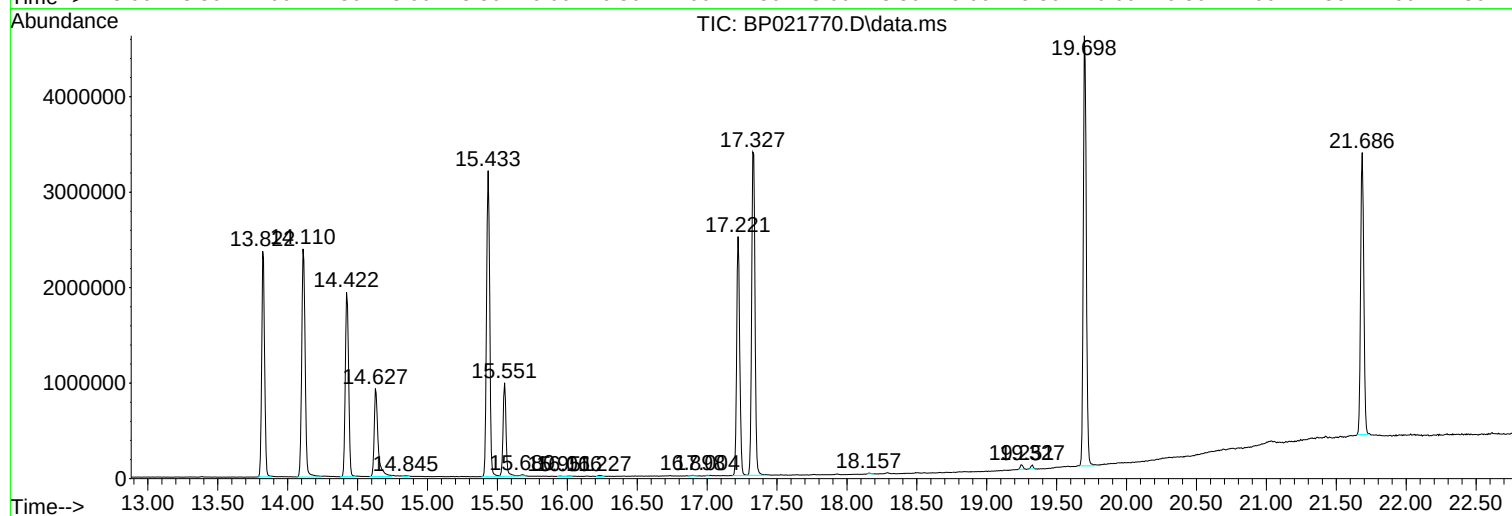
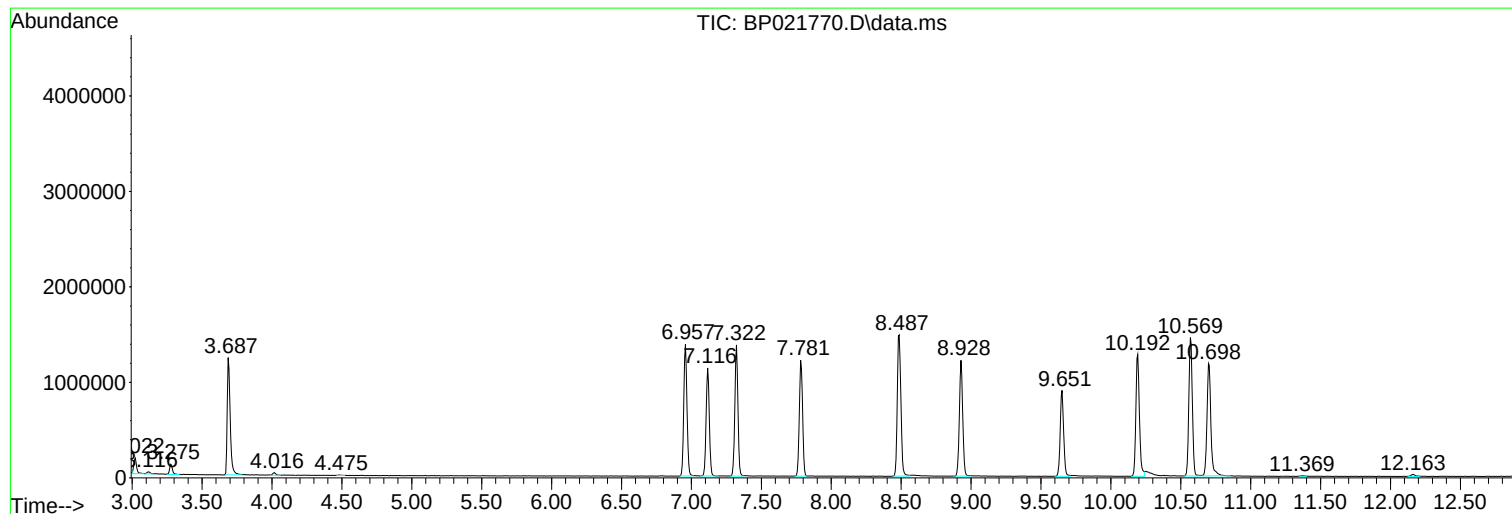
Sum of corrected areas: 75838015

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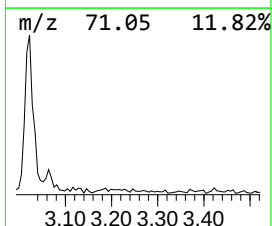
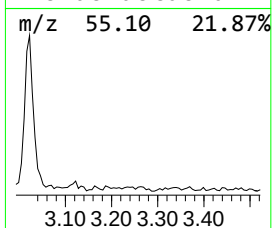
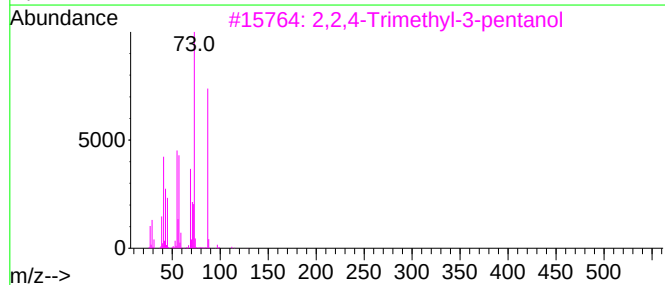
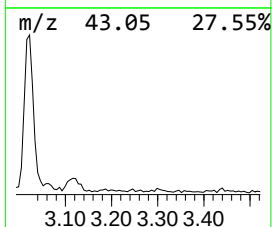
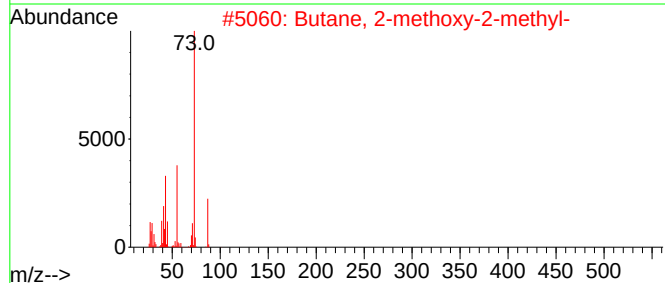
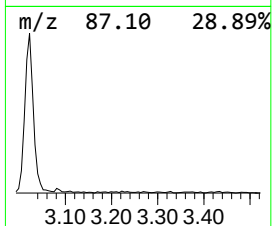
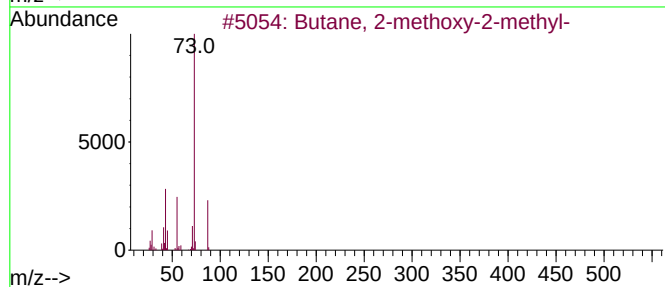
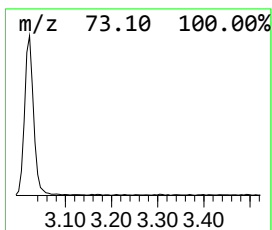
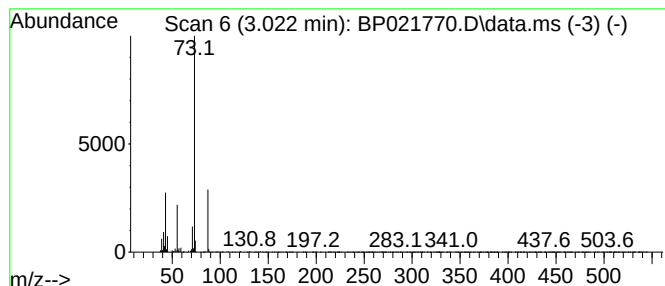
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

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.
3.022	2.22 ng/ul	217955	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	86
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-(2-Methoxyethoxy)-2-methyl-2-p...	162	C8H18O3	1000364-24-4	23
5			N-Ethylformamide	73	C3H7NO	000627-45-2	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...	3.022	2.2	ng/u1	217955	1	7.781	1966530	20.0