

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082624\
 Data File : BP021642.D
 Acq On : 26 Aug 2024 17:21
 Operator : RC/JU
 Sample : P3336-16
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 F7E44

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: BP021642.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	22	rVB	2617318	3525569	73.75%	5.839%
2	3.281	46	50	59	rVB	58636	75851	1.59%	0.126%
3	3.546	91	95	108	rVB	41713	69193	1.45%	0.115%
4	3.693	115	120	135	rBV	703917	1017835	21.29%	1.686%
5	4.028	172	177	181	rVB	20947	28943	0.61%	0.048%
6	4.381	233	237	242	rVB3	15179	21905	0.46%	0.036%
7	4.934	327	331	335	rVB3	7069	10981	0.23%	0.018%
8	5.146	363	367	371	rBV7	3704	6073	0.13%	0.010%
9	5.899	490	495	499	rBV2	16457	25030	0.52%	0.041%
10	6.069	519	524	527	rBV7	3536	6970	0.15%	0.012%
11	6.193	541	545	550	rBV7	2674	5814	0.12%	0.010%
12	6.746	634	639	641	rBV5	3288	5343	0.11%	0.009%
13	6.799	643	648	655	rVB5	4365	11100	0.23%	0.018%
14	6.963	667	676	688	rBV	830426	1372123	28.70%	2.272%
15	7.128	694	704	714	rVB	704057	1144570	23.94%	1.896%
16	7.334	730	739	747	rBV	787320	1318191	27.58%	2.183%
17	7.399	747	750	758	rVV2	16249	35134	0.73%	0.058%
18	7.487	762	765	771	rVB8	3563	5430	0.11%	0.009%
19	7.793	810	817	824	rBV	1068699	1688860	35.33%	2.797%
20	7.857	824	828	832	rVB4	13842	22454	0.47%	0.037%
21	8.487	928	935	950	rBV	986941	1685666	35.26%	2.792%
22	8.940	1004	1012	1027	rBV	766098	1260314	26.36%	2.087%
23	9.110	1036	1041	1046	rVB6	6925	11858	0.25%	0.020%
24	9.387	1079	1088	1093	rVB4	9315	17947	0.38%	0.030%
25	9.498	1101	1107	1114	rBV	60863	103008	2.15%	0.171%
26	9.657	1126	1134	1142	rBV	555472	888053	18.58%	1.471%
27	10.204	1218	1227	1248	rBV	842859	1569712	32.84%	2.600%
28	10.581	1281	1291	1304	rBV	1302292	2302612	48.17%	3.813%
29	10.704	1304	1312	1327	rBV	813676	1562033	32.68%	2.587%
30	11.122	1377	1383	1388	rBV7	13208	27100	0.57%	0.045%
31	11.322	1411	1417	1434	rBV2	74717	201985	4.23%	0.335%
32	11.528	1444	1452	1461	rVB2	9113	20201	0.42%	0.033%
33	12.010	1529	1534	1540	rVB9	7150	12755	0.27%	0.021%
34	12.081	1543	1546	1552	rVB7	4651	8369	0.18%	0.014%
35	12.175	1556	1562	1569	rVB2	17117	30983	0.65%	0.051%
36	12.875	1679	1681	1686	rVB5	3601	6167	0.13%	0.010%

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

37	13.051	1706	1711	1712	rBV5	3371	5155	0.11%	0.009%
38	13.075	1712	1715	1719	rVB6	3551	5702	0.12%	0.009%
39	13.269	1745	1748	1753	rVB6	9689	12741	0.27%	0.021%
40	13.404	1768	1771	1778	rVB8	4987	8017	0.17%	0.013%
41	13.639	1804	1811	1815	rBV10	4683	10545	0.22%	0.017%
42	13.845	1837	1846	1860	rBV	1450143	2256033	47.19%	3.736%
43	14.134	1884	1895	1905	rBV2	1711342	2744366	57.41%	4.545%
44	14.363	1929	1934	1939	rVB7	6748	10432	0.22%	0.017%
45	14.445	1939	1948	1957	rBV	1964222	2942751	61.56%	4.873%
46	14.639	1974	1981	2000	rBV3	543398	1241728	25.98%	2.056%
47	14.869	2014	2020	2027	rVB3	14688	32158	0.67%	0.053%
48	15.198	2072	2076	2080	rBV7	3317	5589	0.12%	0.009%
49	15.275	2084	2089	2091	rVV6	7970	10331	0.22%	0.017%
50	15.328	2095	2098	2104	rBV8	5313	9245	0.19%	0.015%
51	15.457	2111	2120	2127	rBV	2273976	3419100	71.52%	5.662%
52	15.575	2133	2140	2148	rBV	456838	710123	14.86%	1.176%
53	15.704	2158	2162	2168	rVB6	9976	15900	0.33%	0.026%
54	15.916	2196	2198	2203	rVB5	6037	6357	0.13%	0.011%
55	16.028	2213	2217	2221	rBV6	9627	13428	0.28%	0.022%
56	16.210	2242	2248	2252	rBV8	9418	18542	0.39%	0.031%
57	16.645	2319	2322	2325	rBV5	6172	7995	0.17%	0.013%
58	16.798	2346	2348	2354	rVB7	6710	7496	0.16%	0.012%
59	17.022	2382	2386	2389	rBV5	12859	18152	0.38%	0.030%
60	17.245	2417	2424	2434	rVV	2844823	3585049	75.00%	5.937%
61	17.339	2434	2440	2454	rVV2	2295480	3714650	77.71%	6.152%
62	17.445	2454	2458	2465	rVB7	19584	30343	0.63%	0.050%
63	17.657	2490	2494	2500	rVB7	18883	30401	0.64%	0.050%
64	17.839	2522	2525	2530	rBV	8755	10513	0.22%	0.017%
65	17.957	2541	2545	2550	rVB2	24229	33606	0.70%	0.056%
66	18.186	2579	2584	2590	rBV	214741	297890	6.23%	0.493%
67	19.080	2733	2736	2743	rVB8	36899	65231	1.36%	0.108%
68	19.174	2749	2752	2755	rBV4	28874	37153	0.78%	0.062%
69	19.204	2755	2757	2761	rVB	26083	28465	0.60%	0.047%
70	19.269	2765	2768	2772	rVB3	32982	42969	0.90%	0.071%
71	19.333	2776	2779	2784	rVV	36171	67710	1.42%	0.112%
72	19.386	2786	2788	2799	rVB7	52183	104045	2.18%	0.172%
73	19.510	2805	2809	2816	rBV	97584	137393	2.87%	0.228%
74	19.710	2838	2843	2854	rVB	3589012	4639980	97.06%	7.684%
75	21.374	3122	3126	3132	rBV2	146615	283617	5.93%	0.470%
76	21.704	3175	3182	3190	rBV2	2323162	4338693	90.76%	7.185%

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77	24.880	3713	3722	3735	rVB	1745502	4780309	100.00%	7.917%
78	25.115	3754	3762	3775	rVB	1629696	4537113	94.91%	7.514%

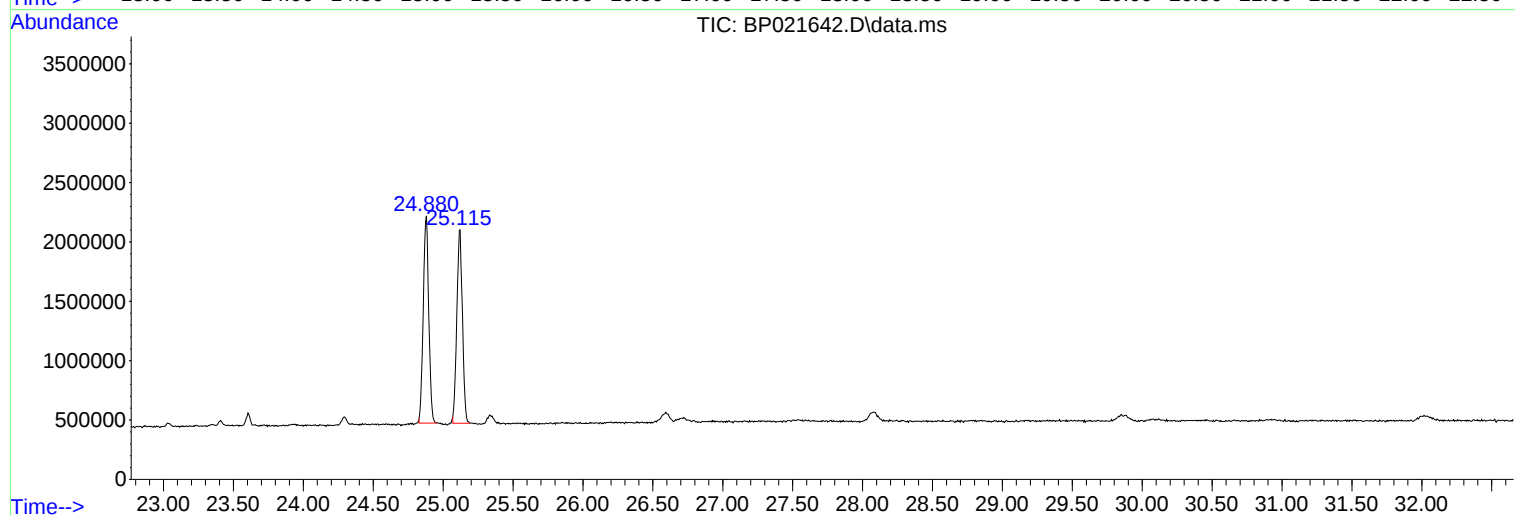
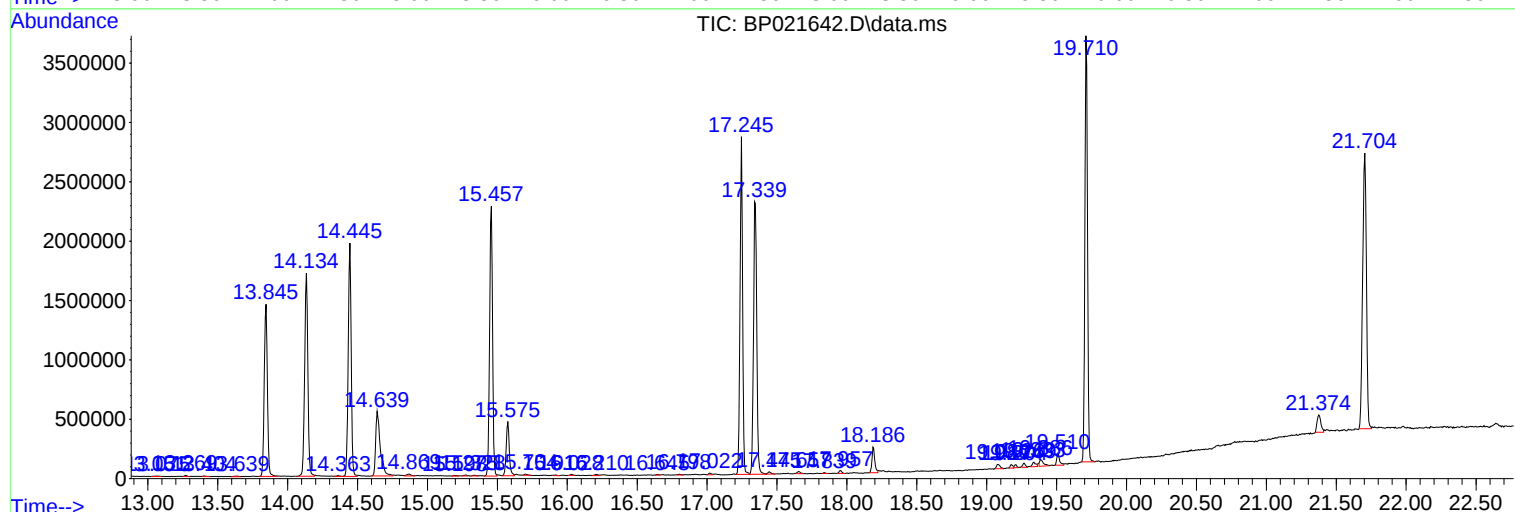
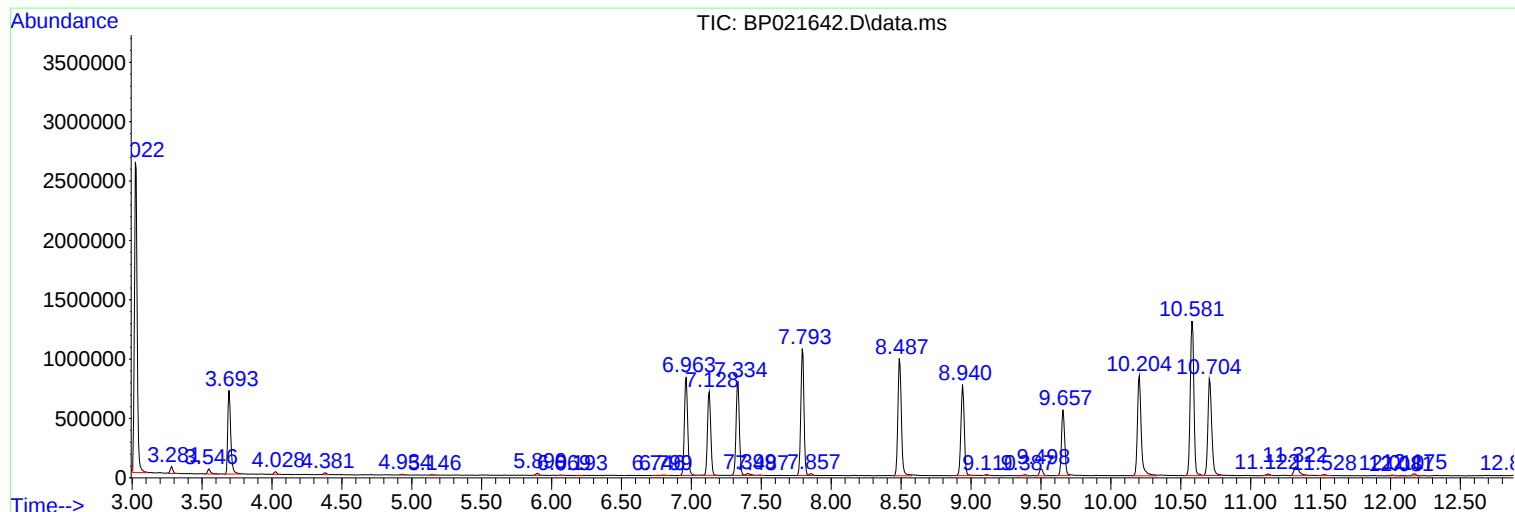
Sum of corrected areas: 60383148

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Instrument :
BNA_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



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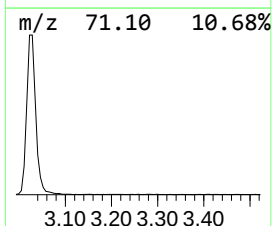
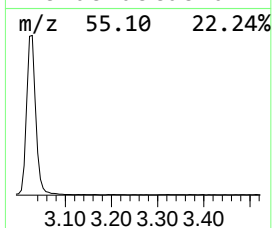
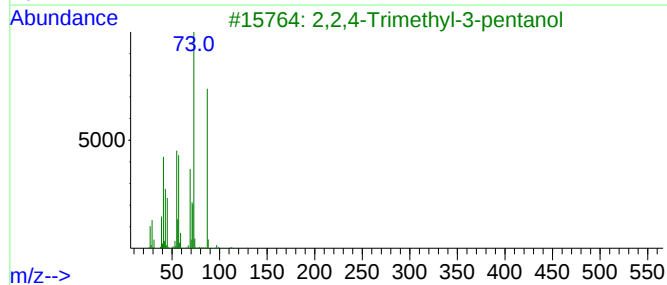
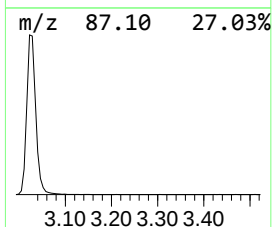
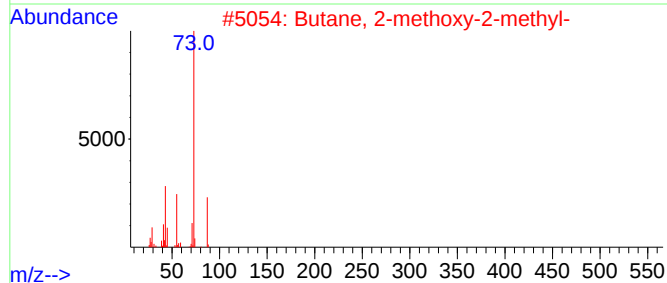
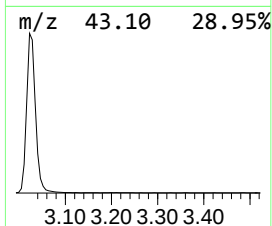
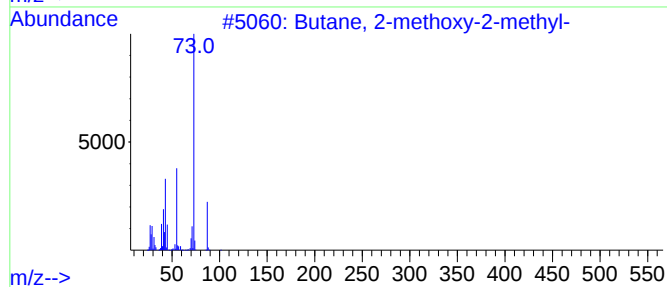
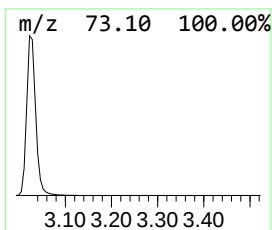
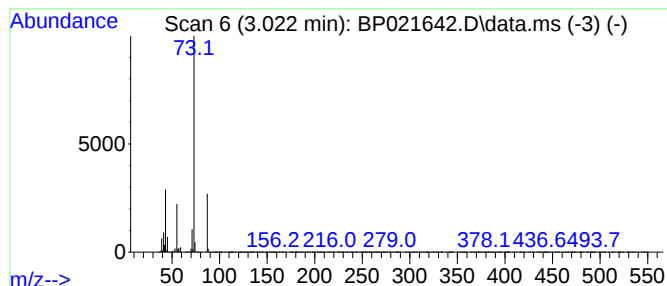
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

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	41.75 ng/ul	3525570	1,4-Dichlorobenzene-d4	7.793

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			Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	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...	3.022	41.8	ng/u1	3525570	1	7.793	1688860	20.0