

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032923\
 Data File : BP014378.D
 Acq On : 29 Mar 2023 23:48
 Operator : CG/JU
 Sample : 02041-02
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 CB956

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-BP032823.M
 Title : SVOA CALIBRATION

Signal : TIC: BP014378.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.393	31	35	38	rBV	35734	46523	0.81%	0.091%
2	3.428	38	41	50	rVB	37608	58531	1.02%	0.114%
3	3.670	77	82	90	rBV	106234	141010	2.45%	0.275%
4	3.828	107	109	124	rBV	167883	288125	5.00%	0.561%
5	3.940	124	128	133	rVV3	16864	30157	0.52%	0.059%
6	3.999	134	138	141	rVV5	8471	13388	0.23%	0.026%
7	4.046	141	146	151	rVB	15780	24265	0.42%	0.047%
8	4.146	159	163	170	rVB	12742	20422	0.35%	0.040%
9	4.258	178	182	188	rBV8	8566	12931	0.22%	0.025%
10	4.534	225	229	233	rVB7	7900	10964	0.19%	0.021%
11	4.593	233	239	248	rBV2	46410	73578	1.28%	0.143%
12	4.723	257	261	269	rVB2	4464	10892	0.19%	0.021%
13	4.958	298	301	302	rVV2	9368	10230	0.18%	0.020%
14	4.981	303	305	310	rVB3	11474	13626	0.24%	0.027%
15	5.081	318	322	329	rBV3	11081	23416	0.41%	0.046%
16	5.705	423	428	436	rBV3	3657	8546	0.15%	0.017%
17	6.087	485	493	498	rBV4	10745	19210	0.33%	0.037%
18	6.393	540	545	552	rVB5	8083	16891	0.29%	0.033%
19	7.117	663	668	672	rBV6	3760	5995	0.10%	0.012%
20	7.181	672	679	690	rBV	129070	258734	4.49%	0.504%
21	7.340	700	706	717	rBV	621334	1016077	17.64%	1.979%
22	7.546	733	741	759	rBV	596360	992739	17.24%	1.934%
23	8.016	814	821	830	rBV	646289	1062123	18.44%	2.069%
24	8.734	935	943	962	rBV	484814	868990	15.09%	1.693%
25	9.034	988	994	997	rBV7	4010	6757	0.12%	0.013%
26	9.193	1014	1021	1035	rBV	804571	1366595	23.73%	2.662%
27	9.569	1076	1085	1090	rBV3	19381	39876	0.69%	0.078%
28	9.922	1138	1145	1158	rBV	685344	1169185	20.30%	2.278%
29	10.387	1222	1224	1230	rVB7	3640	5928	0.10%	0.012%
30	10.463	1230	1237	1258	rBV	869288	1695562	29.44%	3.303%
31	10.840	1293	1301	1309	rBV	920447	1575991	27.37%	3.070%
32	10.987	1318	1326	1349	rBV	725841	1421332	24.68%	2.769%
33	11.310	1375	1381	1385	rVB9	2825	6062	0.11%	0.012%
34	12.181	1523	1529	1535	rBV3	13325	24730	0.43%	0.048%
35	12.240	1535	1539	1546	rVB7	7622	14162	0.25%	0.028%
36	12.369	1554	1561	1567	rBV	38417	78795	1.37%	0.153%

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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-BP032823.M
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37	12.434	1568	1572	1579	rVV2	15290	32751	0.57%	0.064%
38	12.663	1606	1611	1615	rVB8	5594	8943	0.16%	0.017%
39	13.134	1686	1691	1693	rBV5	4547	6489	0.11%	0.013%
40	13.475	1743	1749	1755	rBV2	14497	26062	0.45%	0.051%
41	13.551	1758	1762	1769	rVB5	14668	24873	0.43%	0.048%
42	13.793	1799	1803	1809	rVB7	18165	27417	0.48%	0.053%
43	13.993	1833	1837	1842	rBV8	3349	8269	0.14%	0.016%
44	14.075	1843	1851	1866	rVV	2193750	3102824	53.88%	6.044%
45	14.187	1868	1870	1874	rVB5	6542	8523	0.15%	0.017%
46	14.369	1894	1901	1910	rBV	2173068	3246684	56.38%	6.325%
47	14.593	1934	1939	1946	rVV4	27410	40624	0.71%	0.079%
48	14.675	1946	1953	1962	rVB	1480586	2185361	37.95%	4.257%
49	14.916	1989	1994	2010	rBV2	60069	190263	3.30%	0.371%
50	15.410	2074	2078	2083	rBV2	21507	32804	0.57%	0.064%
51	15.669	2115	2122	2132	rVV	2925228	4192462	72.80%	8.167%
52	15.793	2137	2143	2154	rBV	1007027	1491299	25.90%	2.905%
53	16.034	2179	2184	2193	rVB	117860	163771	2.84%	0.319%
54	16.187	2204	2210	2215	rBV	131920	195468	3.39%	0.381%
55	16.522	2265	2267	2270	rBV3	6243	6130	0.11%	0.012%
56	16.704	2289	2298	2307	rBV2	78843	189110	3.28%	0.368%
57	17.240	2385	2389	2395	rVB2	20776	32328	0.56%	0.063%
58	17.434	2416	2422	2432	rVB	1944871	2669616	46.36%	5.200%
59	17.534	2433	2439	2448	rBV	3530175	4917587	85.39%	9.579%
60	18.298	2565	2569	2578	rBV2	123762	214566	3.73%	0.418%
61	18.604	2619	2621	2626	rVB6	21449	25978	0.45%	0.051%
62	19.339	2743	2746	2750	rVB2	45232	58631	1.02%	0.114%
63	19.410	2754	2758	2763	rVB	45689	70985	1.23%	0.138%
64	19.528	2775	2778	2784	rBV2	62436	91363	1.59%	0.178%
65	19.763	2812	2818	2823	rBV	4427186	5759005	100.00%	11.219%
66	21.198	3059	3062	3071	rBV2	116116	199229	3.46%	0.388%
67	21.522	3112	3117	3123	rBV	1948566	2605095	45.24%	5.075%
68	23.880	3510	3518	3530	rBV2	2302237	4770200	82.83%	9.292%
69	24.039	3539	3545	3557	rVB2	1079481	2307501	40.07%	4.495%

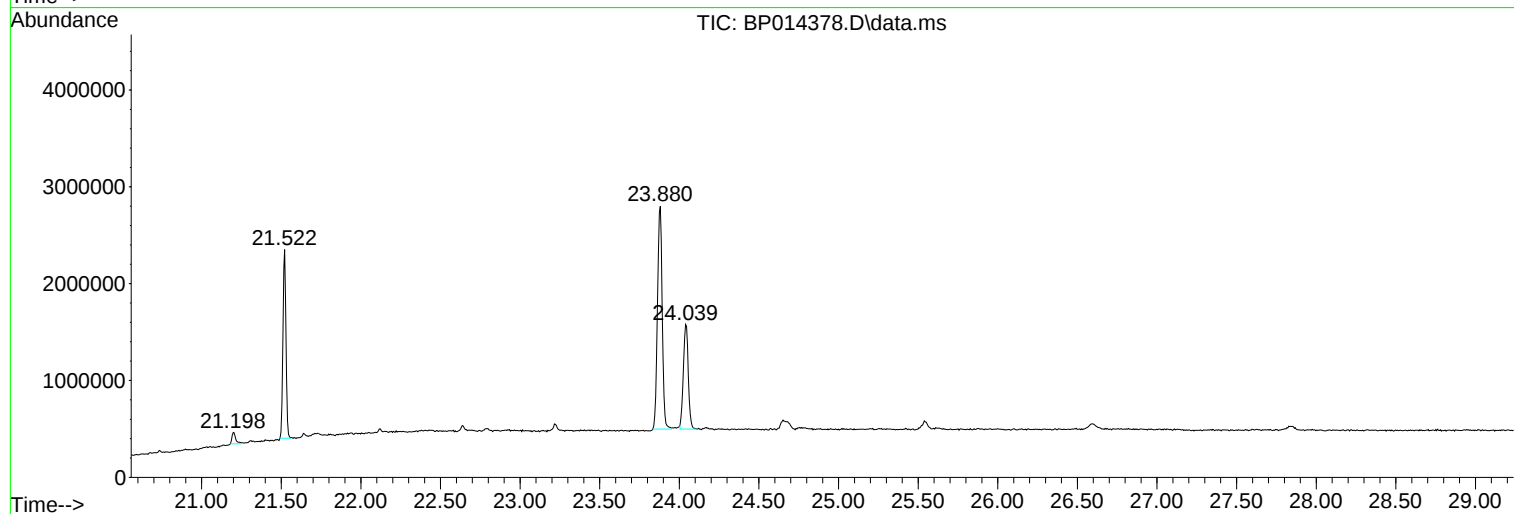
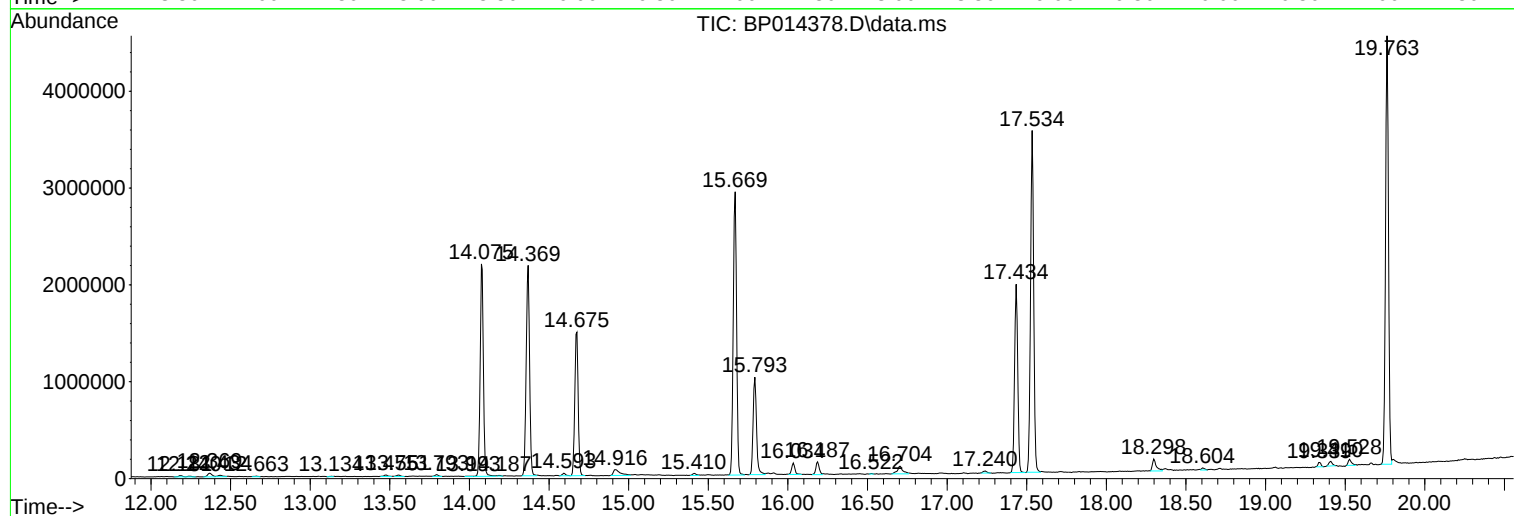
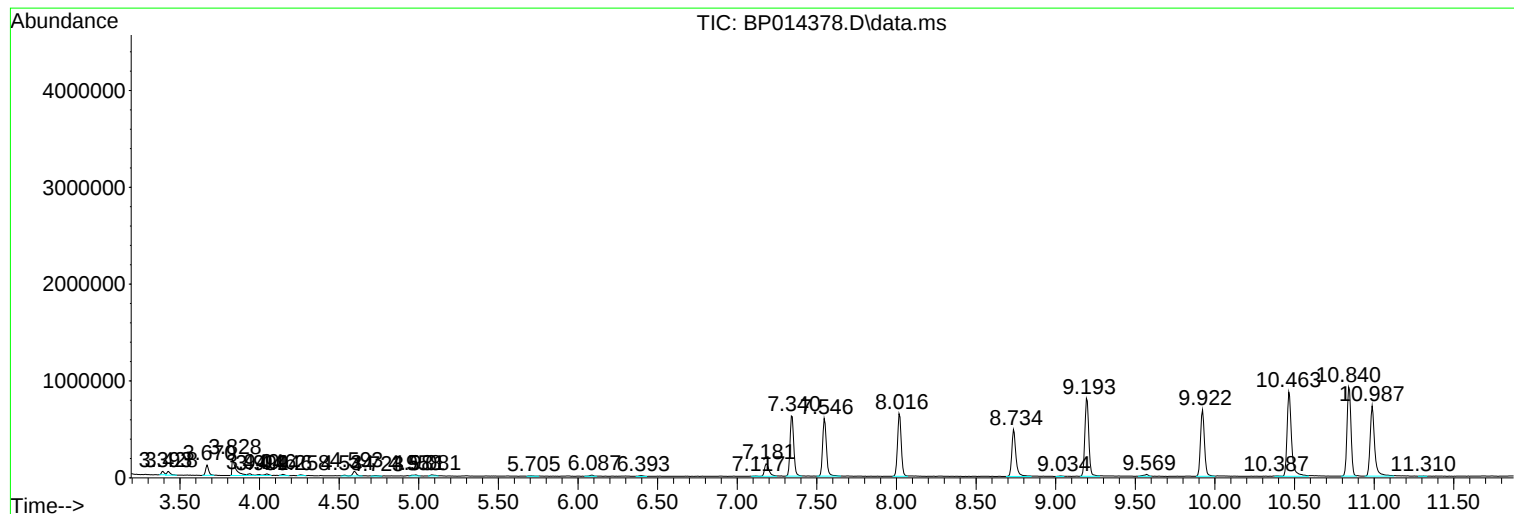
Sum of corrected areas: 51334549

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

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



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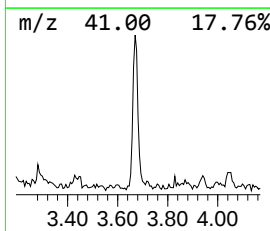
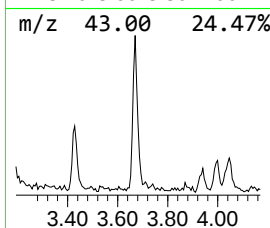
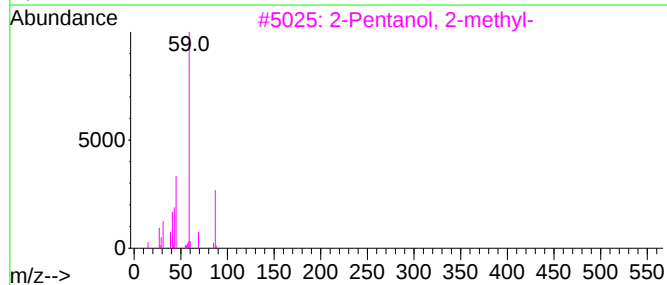
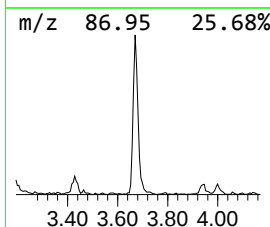
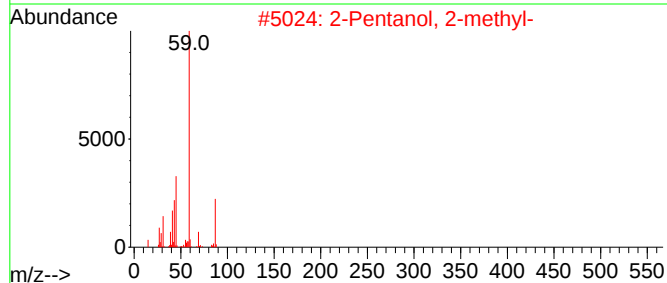
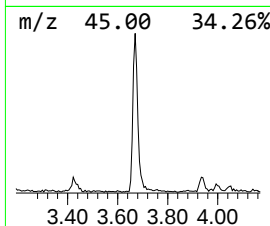
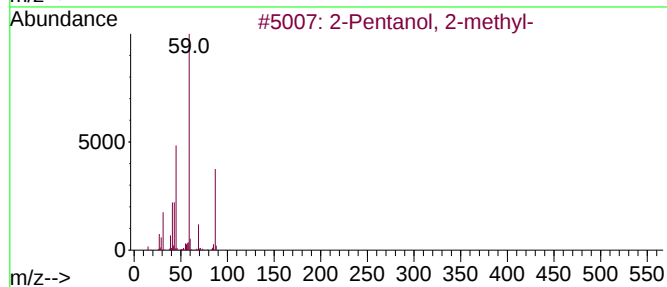
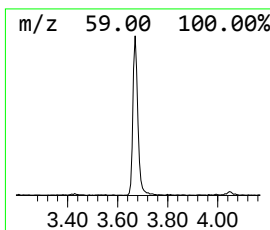
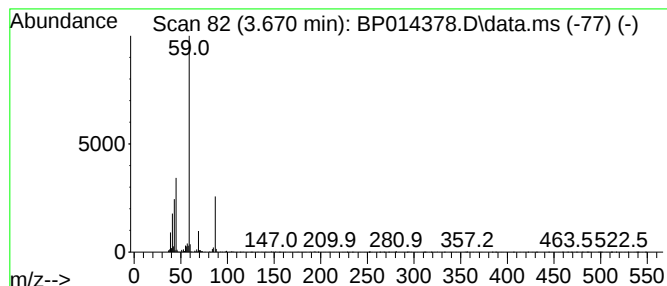
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 2-Pentanol, 2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.670	2.66 ng/ul	141010	1,4-Dichlorobenzene-d4	8.016

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanol, 2-methyl-			102	C6H14O	000590-36-3	90
2	2-Pentanol, 2-methyl-			102	C6H14O	000590-36-3	83
3	2-Pentanol, 2-methyl-			102	C6H14O	000590-36-3	83
4	2-Butanol, 2,3-dimethyl-			102	C6H14O	000594-60-5	64
5	2-Butanol, 2,3-dimethyl-			102	C6H14O	000594-60-5	64



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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
2-Pentanol, 2-m...	3.670	2.7	ng/u1	141010	1	8.016	1062120	20.0