

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103024\
 Data File : BP022810.D
 Acq On : 01 Nov 2024 18:50
 Operator : RC/JU
 Sample : P4627-01
 Misc :
 ALS Vial : 46 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 COCN6

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-BP100724.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP022810.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.175	28	32	38	rVB	16102	20175	0.76%	0.092%
2	3.593	99	103	117	rBV	77551	135841	5.10%	0.621%
3	3.823	138	142	147	rBV7	2912	6017	0.23%	0.028%
4	3.899	151	155	162	rVB4	5376	8011	0.30%	0.037%
5	4.393	234	239	252	rVB	270942	356871	13.39%	1.632%
6	4.793	304	307	313	rVB7	1836	2805	0.11%	0.013%
7	5.193	370	375	377	rBV4	3012	3547	0.13%	0.016%
8	5.317	392	396	405	rVB	9212	20118	0.75%	0.092%
9	5.681	454	458	462	rBV2	4168	5863	0.22%	0.027%
10	6.311	555	565	571	rBV5	8830	18625	0.70%	0.085%
11	6.652	616	623	629	rBV4	6675	15668	0.59%	0.072%
12	6.728	634	636	640	rVV4	3448	5768	0.22%	0.026%
13	6.769	640	643	646	rVV5	3444	4846	0.18%	0.022%
14	6.828	647	653	670	rVV	77978	148584	5.57%	0.680%
15	6.969	670	677	689	rVV	224136	345697	12.97%	1.581%
16	7.169	705	711	722	rBV	254849	455726	17.10%	2.084%
17	7.281	726	730	735	rVB2	5738	9845	0.37%	0.045%
18	7.628	782	789	798	rBV	265477	423467	15.89%	1.937%
19	7.711	798	803	809	rVB7	3885	8187	0.31%	0.037%
20	8.334	902	909	922	rBV	238738	417842	15.68%	1.911%
21	8.446	925	928	936	rVB4	3959	7219	0.27%	0.033%
22	8.769	973	983	998	rBV	302936	524544	19.68%	2.399%
23	8.910	1002	1007	1013	rBV	1590	3817	0.14%	0.017%
24	9.169	1044	1051	1057	rVB4	5415	8290	0.31%	0.038%
25	9.487	1097	1105	1116	rBV	252692	470805	17.66%	2.153%
26	10.016	1188	1195	1214	rBV	355270	711578	26.69%	3.254%
27	10.381	1249	1257	1267	rBV	346739	573956	21.53%	2.625%
28	10.528	1272	1282	1295	rBV	300259	576415	21.62%	2.636%
29	10.934	1344	1351	1357	rBV4	3710	8309	0.31%	0.038%
30	10.987	1357	1360	1364	rVV4	4143	7065	0.27%	0.032%
31	11.046	1364	1370	1382	rVB2	19898	44564	1.67%	0.204%
32	11.452	1434	1439	1441	rBV4	1945	2968	0.11%	0.014%
33	11.599	1457	1464	1469	rBV	20491	34711	1.30%	0.159%
34	11.652	1469	1473	1478	rVB5	4094	7198	0.27%	0.033%
35	11.899	1509	1515	1523	rBV2	13608	26435	0.99%	0.121%
36	11.987	1526	1530	1539	rVB4	4799	8760	0.33%	0.040%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
 Title : SVOA CALIBRATION

37	13.140	1722	1726	1732	rBV7	2278	3818	0.14%	0.017%
38	13.663	1808	1815	1829	rBV	713975	1079067	40.48%	4.935%
39	13.810	1834	1840	1843	rVB8	2887	4461	0.17%	0.020%
40	13.940	1855	1862	1879	rBV2	763266	1174187	44.05%	5.370%
41	14.245	1907	1914	1929	rBV	543088	838180	31.44%	3.833%
42	14.528	1957	1962	1980	rBV3	32487	103555	3.88%	0.474%
43	15.269	2081	2088	2098	rBV	1167377	1586506	59.52%	7.256%
44	15.422	2107	2114	2125	rBV	311561	447869	16.80%	2.048%
45	15.522	2128	2131	2137	rVB6	6854	10975	0.41%	0.050%
46	16.040	2212	2219	2223	rBV8	7072	17653	0.66%	0.081%
47	16.140	2231	2236	2239	rBV2	9397	13548	0.51%	0.062%
48	16.192	2241	2245	2250	rVV4	13177	24154	0.91%	0.110%
49	16.245	2250	2254	2259	rVV2	10489	17925	0.67%	0.082%
50	16.292	2259	2262	2267	rVB4	4638	6120	0.23%	0.028%
51	16.375	2272	2276	2278	rVV5	2223	3564	0.13%	0.016%
52	16.404	2278	2281	2286	rVB6	2191	3266	0.12%	0.015%
53	16.457	2286	2290	2298	rBV9	4875	10679	0.40%	0.049%
54	16.534	2298	2303	2307	rVV2	9463	16314	0.61%	0.075%
55	16.592	2310	2313	2319	rVB6	4713	7454	0.28%	0.034%
56	17.063	2386	2393	2403	rBV	763317	1091992	40.97%	4.994%
57	17.169	2403	2411	2428	rVV	1286040	1882446	70.62%	8.609%
58	17.292	2428	2432	2440	rVB	3624	7857	0.29%	0.036%
59	17.439	2452	2457	2461	rBV7	1780	3660	0.14%	0.017%
60	17.757	2508	2511	2518	rVB10	1970	3463	0.13%	0.016%
61	17.886	2529	2533	2538	rBV6	6195	8972	0.34%	0.041%
62	17.992	2547	2551	2560	rBV3	18711	31971	1.20%	0.146%
63	18.069	2560	2564	2569	rVB3	3891	6618	0.25%	0.030%
64	19.086	2733	2737	2743	rBV2	16856	25922	0.97%	0.119%
65	19.169	2748	2751	2755	rBV2	11785	18701	0.70%	0.086%
66	19.333	2774	2779	2786	rBV7	26319	43487	1.63%	0.199%
67	19.551	2810	2816	2829	rBV	1562570	2393871	89.80%	10.948%
68	21.510	3143	3149	3160	rVB	903630	1398440	52.46%	6.396%
69	24.539	3654	3664	3681	rVB2	840070	2665646	100.00%	12.191%
70	24.768	3695	3703	3715	rVB	544484	1482837	55.63%	6.782%

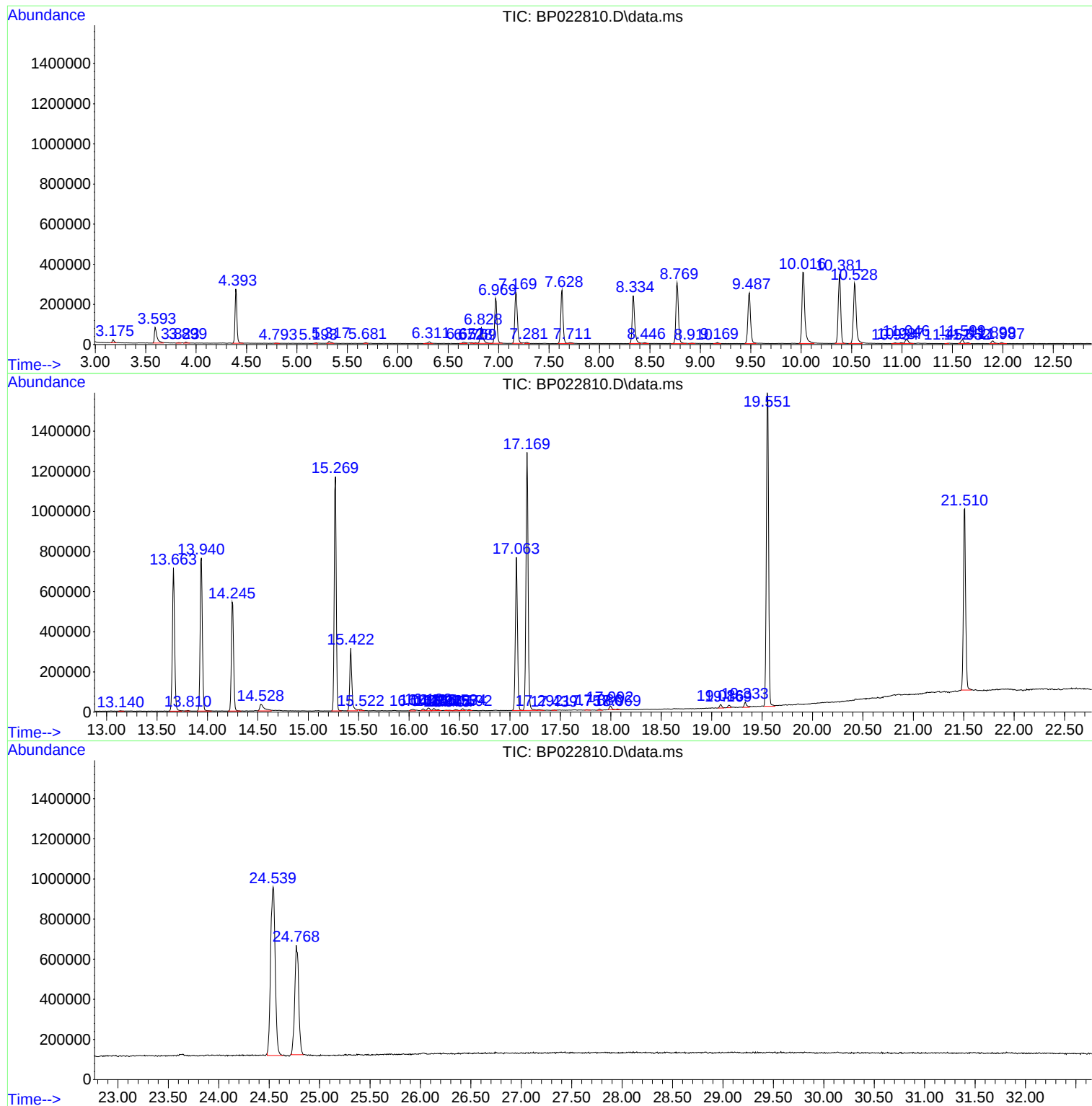
Sum of corrected areas: 21865315

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION

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



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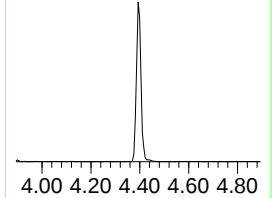
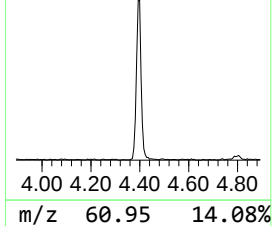
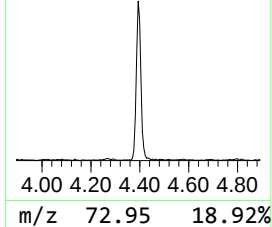
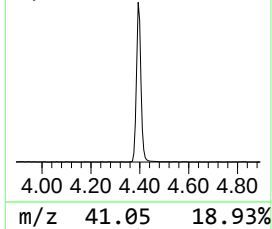
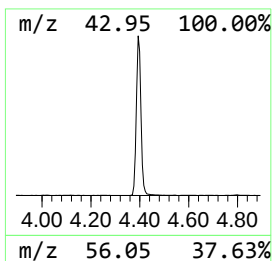
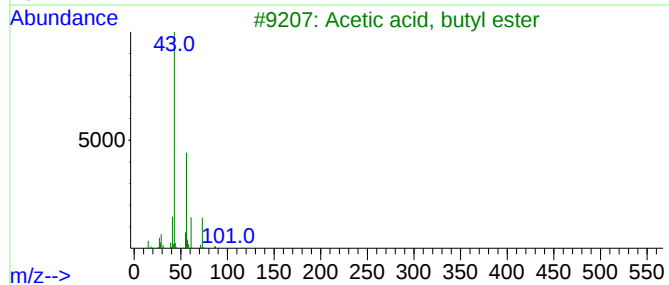
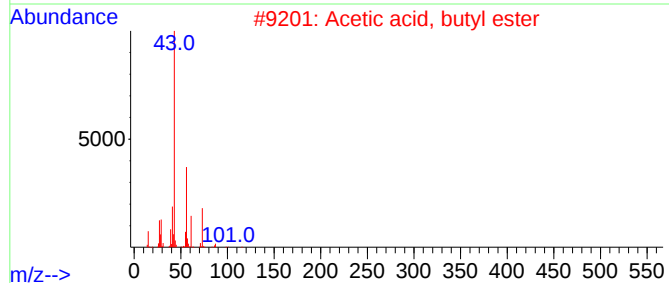
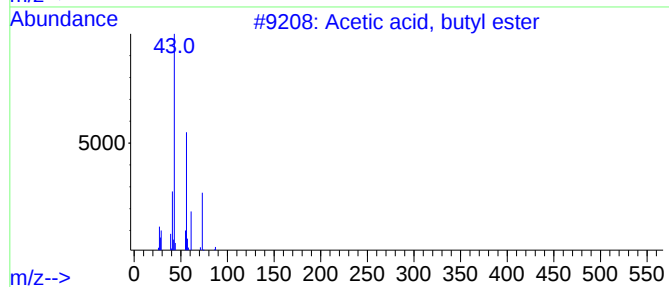
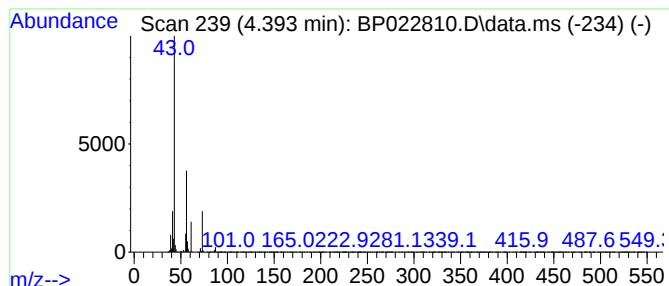
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Acetic acid, butyl ester Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.393	16.85 ng/ul	356871	1,4-Dichlorobenzene-d4	7.628

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetic acid, butyl ester	116	C6H12O2	000123-86-4	90		
2	Acetic acid, butyl ester	116	C6H12O2	000123-86-4	83		
3	Acetic acid, butyl ester	116	C6H12O2	000123-86-4	78		
4	Acetic acid, butyl ester	116	C6H12O2	000123-86-4	64		
5	Isobutyl acetate	116	C6H12O2	000110-19-0	56		



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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
Acetic acid, bu...	4.393	16.9	ng/ul	356871	1	7.628	423467	20.0