

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN102922\
 Data File : BN022364.D
 Acq On : 27 Oct 2022 16:46
 Operator : CG/JU
 Sample : PB148501BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SBLK501

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_N\Methods\SFAM-EPA-BN101222.M
 Title : SVOA CALIBRATION

Signal : TIC: BN022364.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.246	40	44	57	rVB	54849	74296	2.76%	0.220%
2	3.693	117	120	136	rBV	91778	162009	6.03%	0.479%
3	4.022	171	176	180	rVB	13837	19887	0.74%	0.059%
4	5.169	368	371	383	rVB4	2526	7343	0.27%	0.022%
5	7.093	691	698	717	rBV	788073	1251660	46.56%	3.703%
6	7.263	720	727	740	rBV	664231	1023938	38.09%	3.030%
7	7.457	753	760	776	rVB	759459	1234678	45.93%	3.653%
8	7.934	834	841	855	rBV	627639	1001204	37.24%	2.962%
9	8.657	956	964	977	rBV	925051	1501373	55.85%	4.442%
10	9.116	1034	1042	1055	rBV	734466	1188670	44.22%	3.517%
11	9.846	1158	1166	1176	rBV	578357	976492	36.32%	2.889%
12	10.387	1249	1258	1278	rBV	818123	1451765	54.00%	4.295%
13	10.763	1314	1322	1331	rBV	834001	1447458	53.84%	4.283%
14	10.922	1341	1349	1358	rBV	15738	31841	1.18%	0.094%
15	12.363	1586	1594	1599	rBV2	12955	22090	0.82%	0.065%
16	13.581	1797	1801	1805	rVB3	2492	3369	0.13%	0.010%
17	14.022	1870	1876	1887	rBV	1345325	1841670	68.51%	5.449%
18	14.304	1917	1924	1931	rBV	1623841	2339260	87.02%	6.921%
19	14.610	1970	1976	1987	rVB	1171261	1732076	64.43%	5.125%
20	14.816	2004	2011	2032	rBV	569555	880147	32.74%	2.604%
21	15.016	2042	2045	2051	rVB5	2376	3578	0.13%	0.011%
22	15.604	2139	2145	2158	rVB	1885602	2688285	100.00%	7.954%
23	15.734	2161	2167	2182	rBV	515800	707707	26.33%	2.094%
24	15.845	2182	2186	2190	rVB3	7443	10447	0.39%	0.031%
25	15.969	2203	2207	2211	rBV3	3951	5383	0.20%	0.016%
26	16.163	2235	2240	2242	rBV4	1935	2968	0.11%	0.009%
27	16.481	2290	2294	2303	rVB6	5660	10800	0.40%	0.032%
28	17.122	2395	2403	2408	rBV2	19441	36143	1.34%	0.107%
29	17.228	2418	2421	2425	rBV4	4473	6123	0.23%	0.018%
30	17.275	2425	2429	2435	rVB2	18858	26734	0.99%	0.079%
31	17.369	2438	2445	2452	rBV	1177697	1691277	62.91%	5.004%
32	17.469	2455	2462	2477	rVV2	1749329	2598741	96.67%	7.689%
33	17.651	2488	2493	2499	rBV4	7396	14662	0.55%	0.043%
34	17.757	2507	2511	2516	rBV4	4172	6972	0.26%	0.021%
35	17.857	2525	2528	2531	rBV3	4556	5625	0.21%	0.017%
36	18.033	2555	2558	2564	rVB2	15277	19031	0.71%	0.056%

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Peak Location: TOP

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37	18.257	2593	2596	2600	rVB3	4759	5511	0.21%	0.016%
38	18.339	2605	2610	2618	rVV	42607	63015	2.34%	0.186%
39	18.551	2643	2646	2650	rBV4	7920	11235	0.42%	0.033%
40	18.686	2665	2669	2673	rBV2	28173	36437	1.36%	0.108%
41	19.186	2751	2754	2759	rBV3	8115	10905	0.41%	0.032%
42	19.327	2774	2778	2785	rBV3	21033	36507	1.36%	0.108%
43	19.404	2787	2791	2798	rVV	17368	31051	1.16%	0.092%
44	19.769	2847	2853	2863	rBV	1813785	2445051	90.95%	7.234%
45	20.510	2975	2979	2986	rBV	136026	178942	6.66%	0.529%
46	21.474	3140	3143	3149	rBV	44174	49831	1.85%	0.147%
47	21.569	3154	3159	3165	rBV	1077513	1355279	50.41%	4.010%
48	23.868	3543	3550	3563	rVB	1071401	2193729	81.60%	6.491%
49	24.033	3571	3578	3590	rVB	650633	1355363	50.42%	4.010%

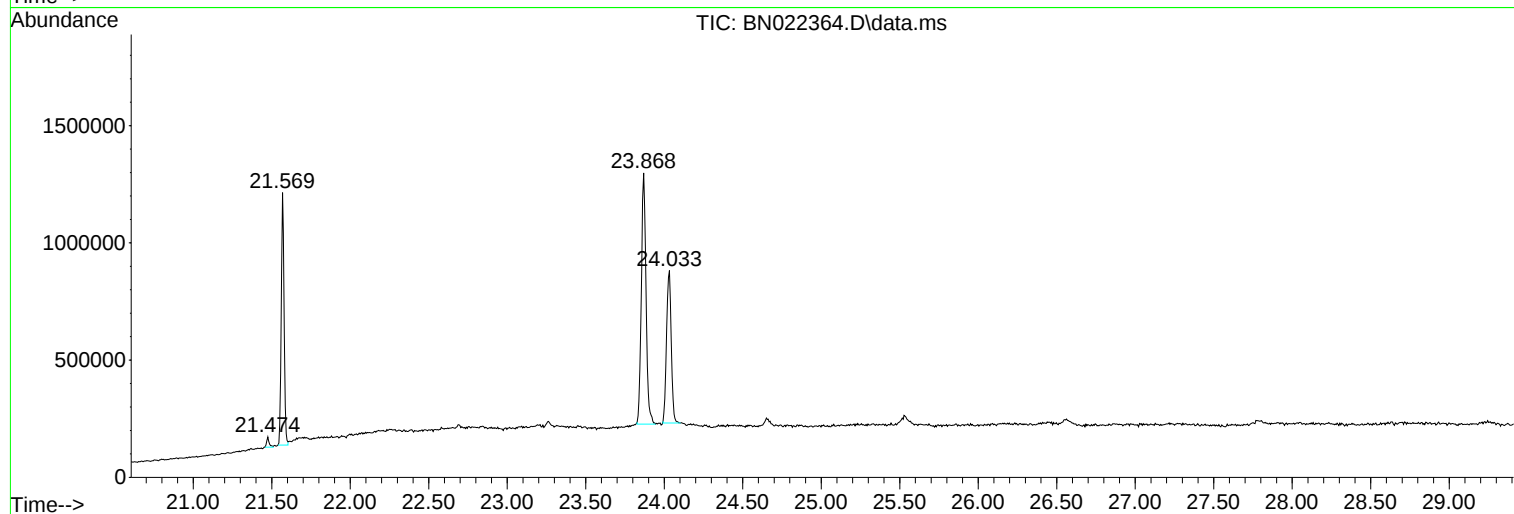
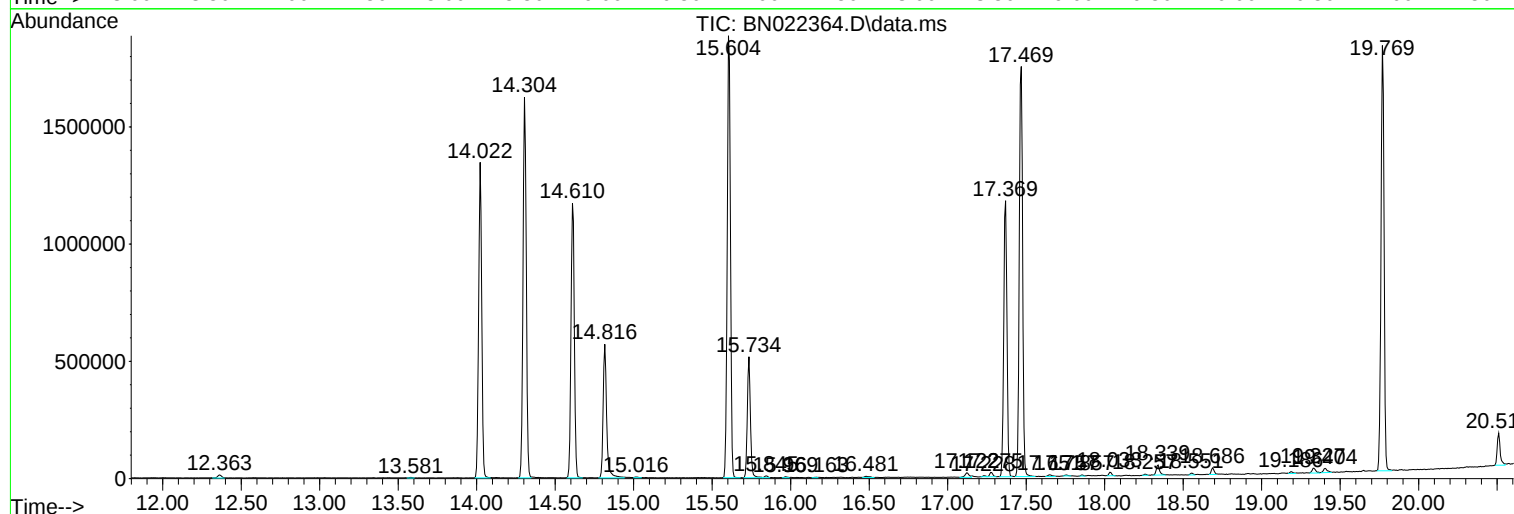
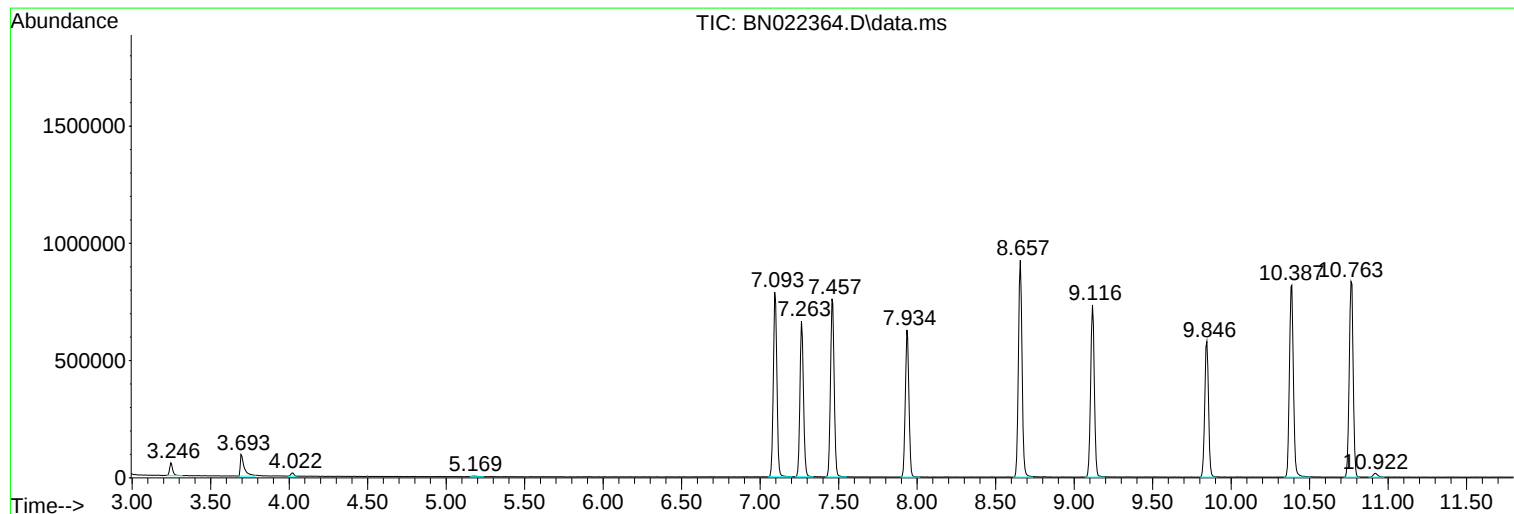
Sum of corrected areas: 33798558

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN101222.M
Quant Title : SVOA CALIBRATION

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



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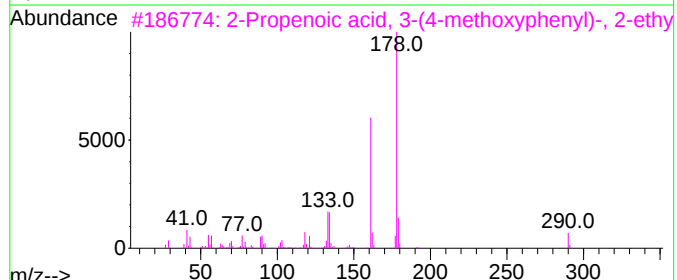
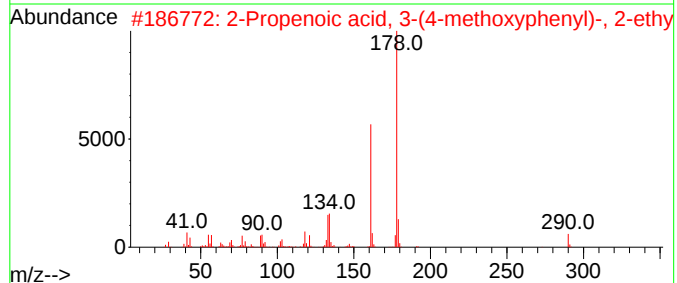
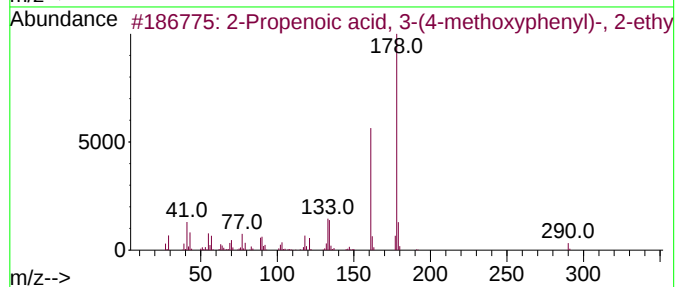
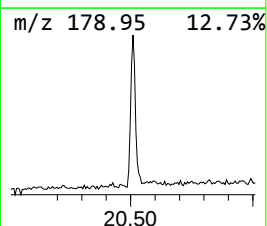
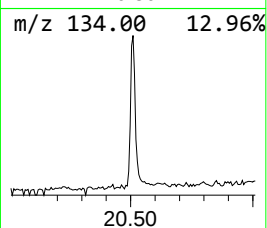
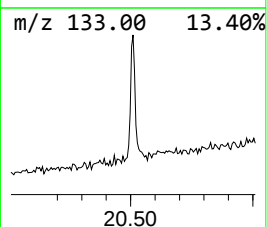
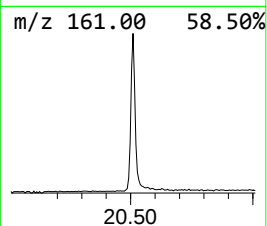
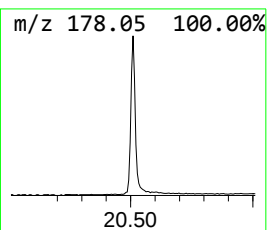
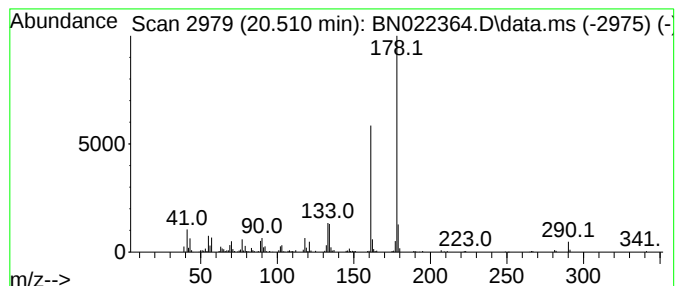
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN101222.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 2-Propenoic acid, 3-(4-meth... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.510	2.64 ng/ul	178942	Chrysene-d12	21.569

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propenoic acid, 3-(4-methoxyph...	290	C18H26O3	005466-77-3	99		
2	2-Propenoic acid, 3-(4-methoxyph...	290	C18H26O3	005466-77-3	99		
3	2-Propenoic acid, 3-(4-methoxyph...	290	C18H26O3	005466-77-3	99		
4	2-Ethylhexyl trans-4-methoxycinn...	290	C18H26O3	083834-59-7	94		
5	2-Propenoic acid, 3-(4-methoxyph...	290	C18H26O3	005466-77-3	93		



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Quant Title : SVOA CALIBRATION

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-Propenoic aci...	20.510	2.6	ng/u1	178942	5	21.569	1355280	20.0