

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081822\
 Data File : BP011486.D
 Acq On : 18 Aug 2022 17:30
 Operator : CG/JU
 Sample : N4221-14
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 COKN6

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-BP081622.M
 Title : SVOA CALIBRATION

Signal : TIC: BP011486.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.034	3	8	16	rBV	5091944	7296961	62.91%	18.479%
2	3.099	16	19	34	rVB	744911	997414	8.60%	2.526%
3	3.493	84	86	96	rBV	42502	67402	0.58%	0.171%
4	3.593	99	103	107	rVB	9977	13264	0.11%	0.034%
5	3.693	115	120	128	rVB	15458	24277	0.21%	0.061%
6	3.781	131	135	141	rBV	14694	20356	0.18%	0.052%
7	3.905	151	156	160	rBV	21890	31297	0.27%	0.079%
8	4.299	216	223	237	rBV	8319250	11598997	100.00%	29.374%
9	4.399	237	240	245	rVB2	14352	18673	0.16%	0.047%
10	5.022	341	346	354	rVB	119571	162372	1.40%	0.411%
11	6.740	633	638	648	rBV	75815	121040	1.04%	0.307%
12	6.904	659	666	680	rBV	299541	460883	3.97%	1.167%
13	7.087	690	697	710	rBV	295439	479426	4.13%	1.214%
14	7.557	770	777	785	rBV	357243	557852	4.81%	1.413%
15	8.275	893	899	908	rBV	218625	347207	2.99%	0.879%
16	8.722	963	975	993	rBV	367226	609289	5.25%	1.543%
17	9.434	1087	1096	1108	rBV	296922	492809	4.25%	1.248%
18	9.969	1177	1187	1197	rBV	384282	646886	5.58%	1.638%
19	10.345	1244	1251	1258	rBV	451940	745729	6.43%	1.889%
20	10.498	1270	1277	1290	rBV	286241	488943	4.22%	1.238%
21	13.069	1708	1714	1720	rBV4	17876	28931	0.25%	0.073%
22	13.657	1806	1814	1825	rBV	731568	1007844	8.69%	2.552%
23	13.875	1843	1851	1854	rBV	342544	510537	4.40%	1.293%
24	13.922	1854	1859	1871	rVB	775145	1172322	10.11%	2.969%
25	14.228	1905	1911	1919	rBV	614587	881400	7.60%	2.232%
26	14.469	1944	1952	1963	rBV2	43478	96137	0.83%	0.243%
27	14.669	1980	1986	1994	rVB8	6289	14042	0.12%	0.036%
28	15.233	2075	2082	2089	rBV	1015034	1377590	11.88%	3.489%
29	15.369	2096	2105	2119	rBV	405061	543478	4.69%	1.376%
30	17.004	2376	2383	2391	rBV	689940	988367	8.52%	2.503%
31	17.104	2393	2400	2412	rVV	1174060	1641780	14.15%	4.158%
32	17.704	2496	2502	2506	rBV3	10971	16134	0.14%	0.041%
33	17.921	2534	2539	2545	rBV3	17894	26959	0.23%	0.068%
34	18.939	2706	2712	2715	rBV3	15739	24679	0.21%	0.062%
35	19.010	2719	2724	2727	rVV2	16001	21670	0.19%	0.055%
36	19.363	2779	2784	2792	rBV	1334140	1779560	15.34%	4.507%

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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

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37	20.868	3036	3040	3046	rBV5	45658	80157	0.69%	0.203%
38	21.133	3080	3085	3092	rBV	855192	1063948	9.17%	2.694%
39	23.280	3443	3450	3457	rBV	1054455	1899513	16.38%	4.810%
40	23.427	3469	3475	3485	rVB2	583995	1131494	9.76%	2.865%

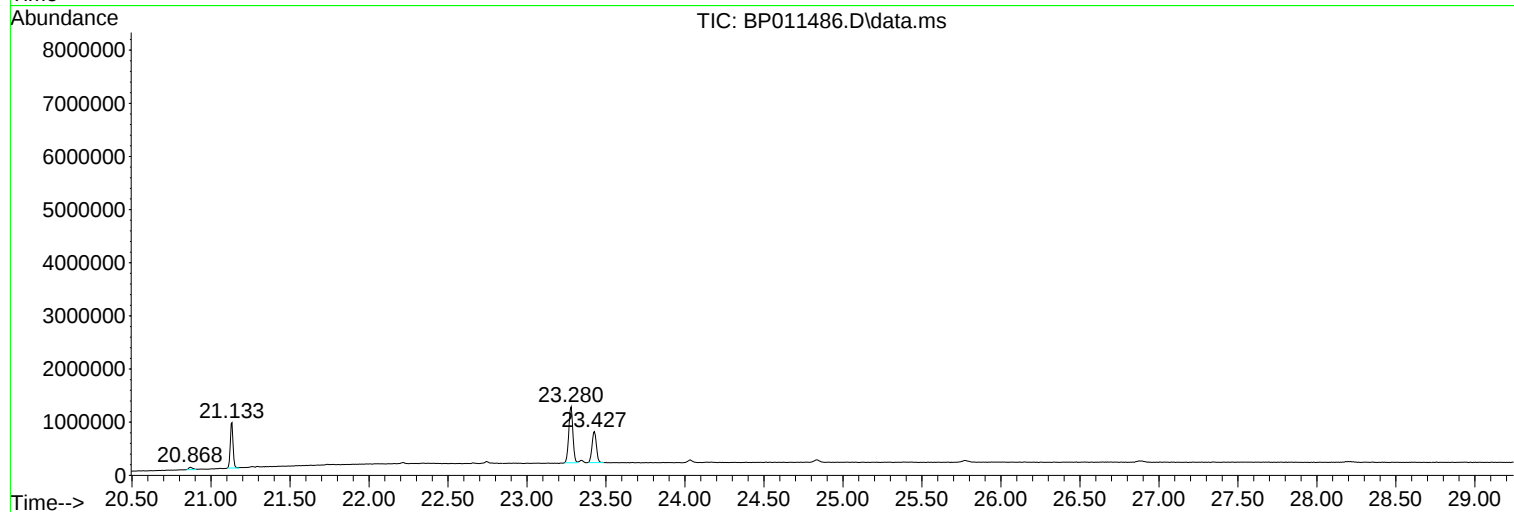
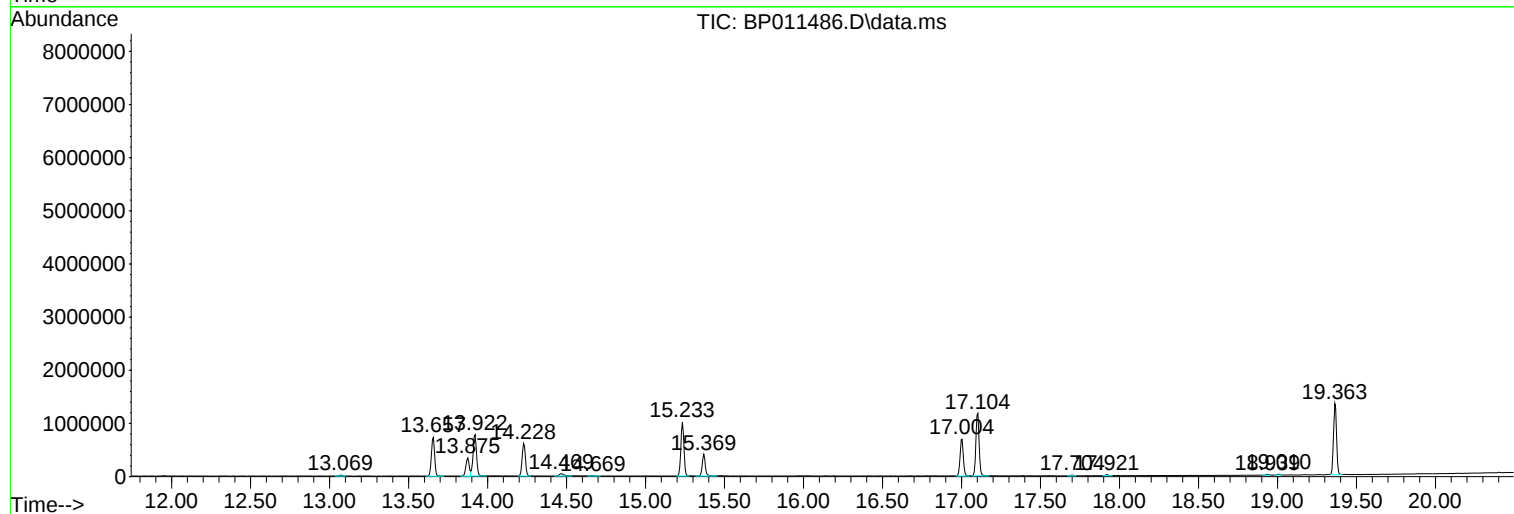
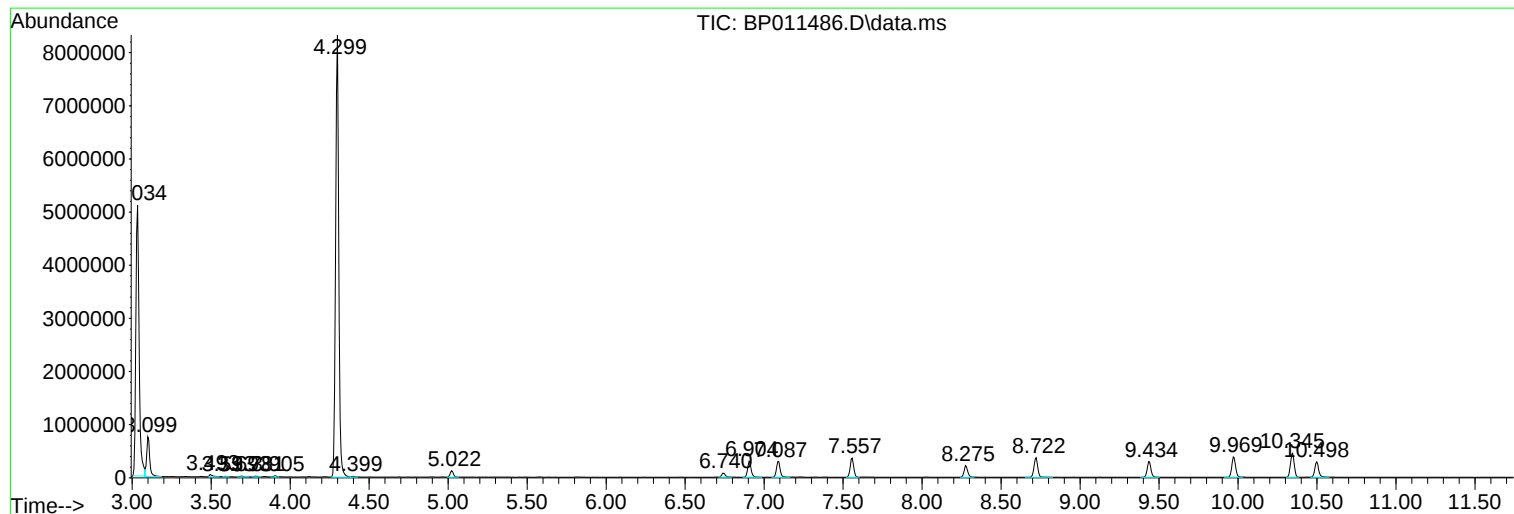
Sum of corrected areas: 39487619

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081622.M
Quant Title : SVOA CALIBRATION

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



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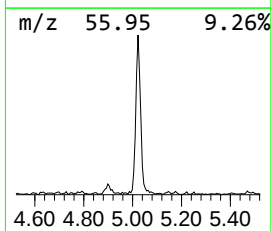
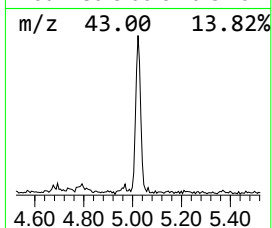
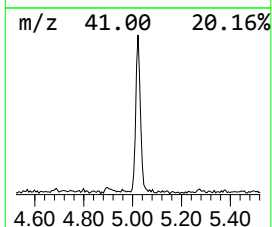
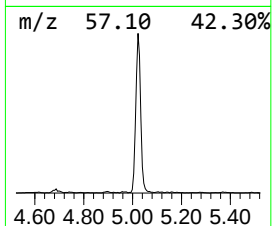
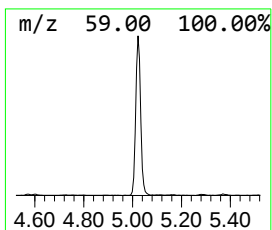
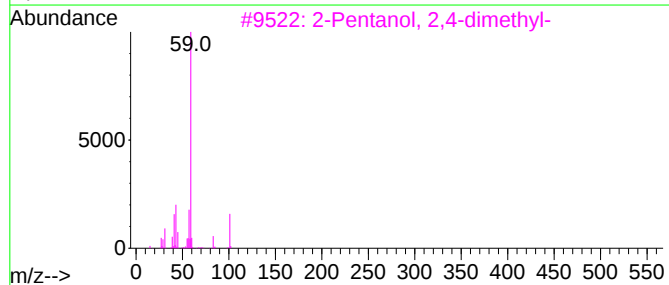
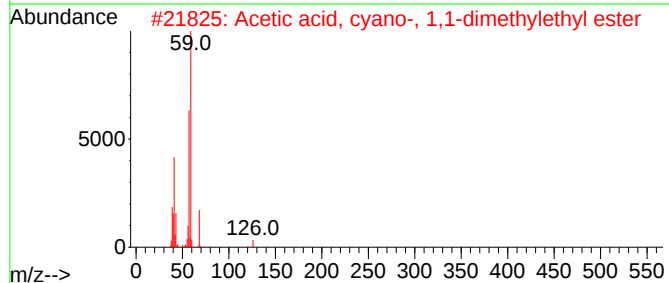
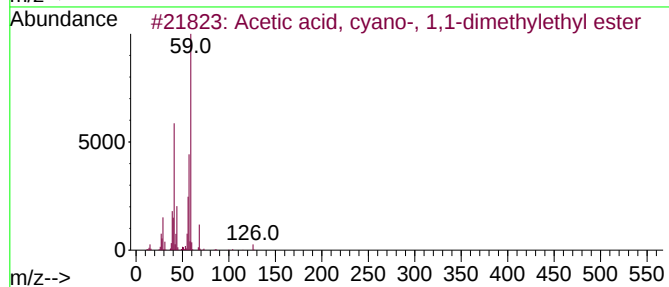
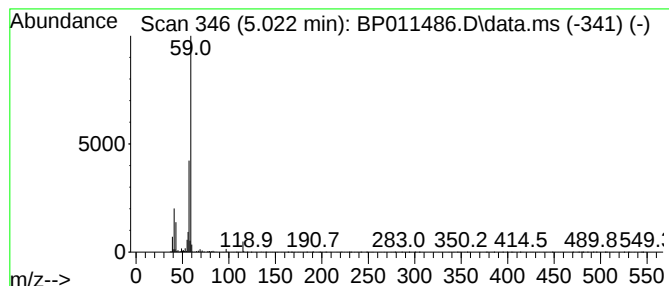
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TIC Integration Parameters: LSCINT.P

Peak Number 2 Acetic acid, cyano-, 1,1-di... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.022	5.82 ng/ul	162372	1,4-Dichlorobenzene-d4	7.557

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	50
2			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	50
3			2-Pentanol, 2,4-dimethyl-	116	C7H16O	000625-06-9	40
4			2-Hexanol, methyl ether	116	C7H16O	1000333-92-0	9
5			Butanamide, 3,3-dimethyl-	115	C6H13NO	000926-04-5	9



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Quant Title : SVOA CALIBRATION

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TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Acetic acid, cy...	5.022	5.8	ng/u1	162372	1	7.557	557852	20.0