

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081822\
 Data File : BP011487.D
 Acq On : 18 Aug 2022 18:05
 Operator : CG/JU
 Sample : N4221-15
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 COKN7

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: BP011487.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	4451749	6668417	52.09%	16.774%
2	3.099	16	19	37	rVB	690681	904567	7.07%	2.275%
3	3.487	82	85	98	rVV	83124	133268	1.04%	0.335%
4	3.693	115	120	127	rVB	15411	24212	0.19%	0.061%
5	3.781	131	135	140	rVB	12954	17523	0.14%	0.044%
6	3.905	150	156	161	rBV	23788	33608	0.26%	0.085%
7	4.299	216	223	236	rBV	9147324	12802355	100.00%	32.203%
8	4.399	236	240	247	rVB	13632	20582	0.16%	0.052%
9	5.022	341	346	354	rVB	130346	171826	1.34%	0.432%
10	6.740	632	638	649	rBV	81744	131466	1.03%	0.331%
11	6.905	657	666	680	rBV	291881	449242	3.51%	1.130%
12	7.087	691	697	710	rBV	307490	475881	3.72%	1.197%
13	7.228	716	721	727	rVB6	8372	17062	0.13%	0.043%
14	7.558	769	777	786	rBV	361492	567277	4.43%	1.427%
15	8.275	893	899	909	rBV	220701	354540	2.77%	0.892%
16	8.722	966	975	983	rBV	353061	576147	4.50%	1.449%
17	9.434	1087	1096	1110	rBV	278949	474698	3.71%	1.194%
18	9.969	1180	1187	1202	rBV	368707	625792	4.89%	1.574%
19	10.340	1243	1250	1262	rVB	454272	755739	5.90%	1.901%
20	10.499	1270	1277	1289	rBV	329155	565333	4.42%	1.422%
21	11.951	1518	1524	1533	rVB2	7651	14934	0.12%	0.038%
22	13.069	1709	1714	1721	rBV2	15135	24975	0.20%	0.063%
23	13.651	1807	1813	1821	rBV	737395	1001697	7.82%	2.520%
24	13.875	1844	1851	1854	rBV	344080	520628	4.07%	1.310%
25	13.922	1854	1859	1873	rVB	721735	1108223	8.66%	2.788%
26	14.228	1905	1911	1919	rBV	621941	882381	6.89%	2.220%
27	14.463	1942	1951	1969	rBV	44454	104000	0.81%	0.262%
28	14.663	1977	1985	1990	rVB	6183	14288	0.11%	0.036%
29	15.234	2074	2082	2089	rBV	981090	1359046	10.62%	3.419%
30	15.369	2097	2105	2118	rBV	399573	525116	4.10%	1.321%
31	16.998	2376	2382	2392	rBV	703677	991804	7.75%	2.495%
32	17.098	2393	2399	2410	rBV2	1136316	1591782	12.43%	4.004%
33	17.922	2534	2539	2543	rBV3	12893	19941	0.16%	0.050%
34	18.939	2708	2712	2716	rBV4	14021	19629	0.15%	0.049%
35	19.004	2720	2723	2728	rVV	12747	17335	0.14%	0.044%
36	19.363	2779	2784	2794	rBV	1375736	1752681	13.69%	4.409%

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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	21.133	3080	3085	3091	rBV	862099	1074845	8.40%	2.704%
38	23.274	3443	3449	3456	rBV2	1020735	1823554	14.24%	4.587%
39	23.427	3468	3475	3483	rVB2	597044	1138797	8.90%	2.865%

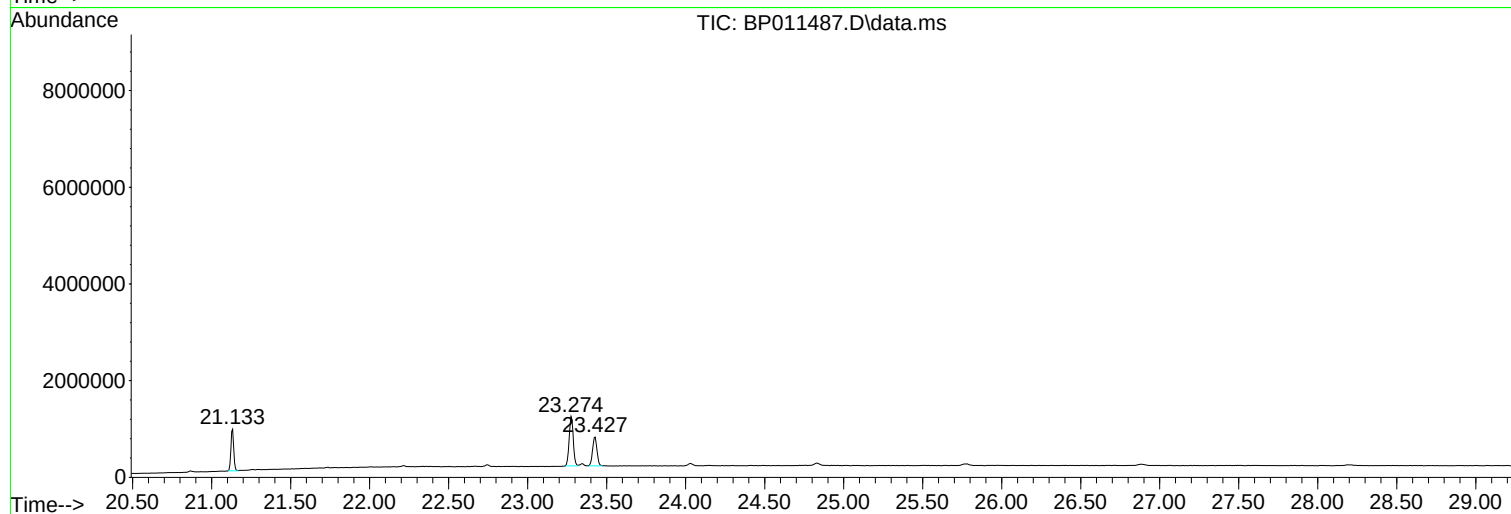
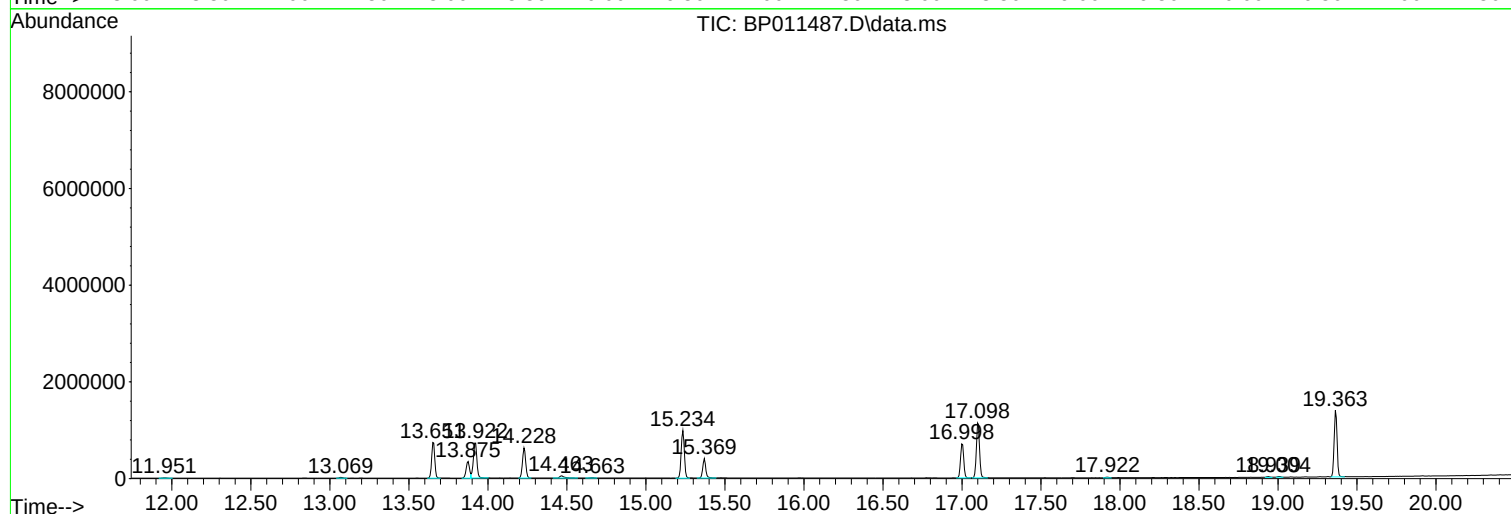
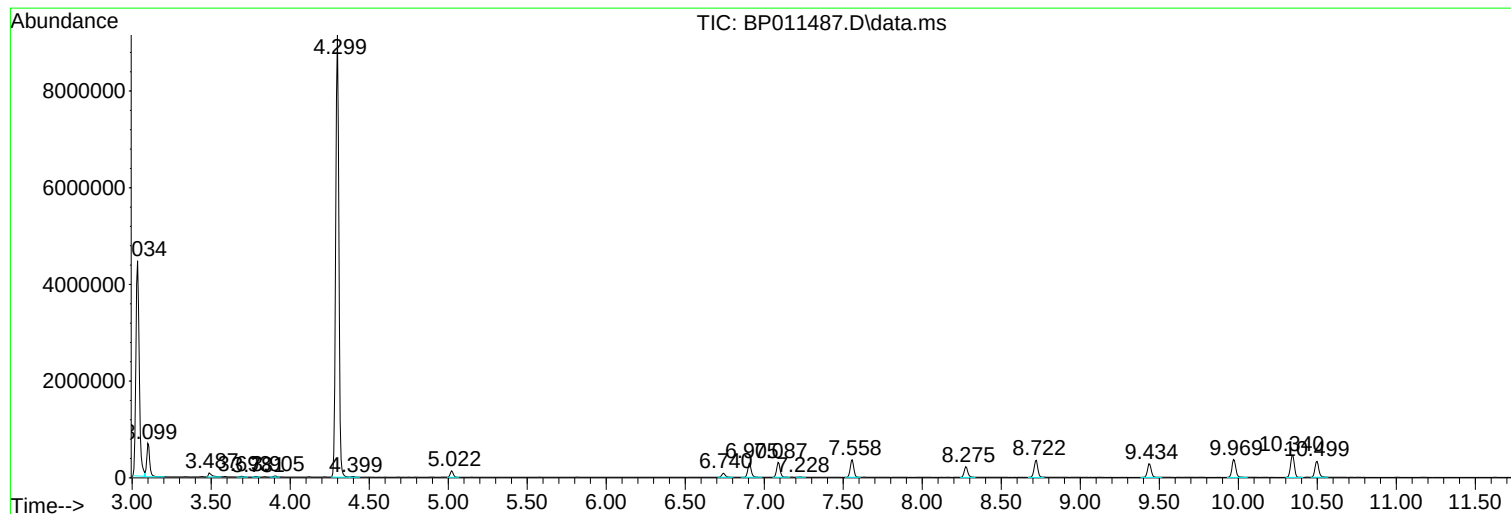
Sum of corrected areas: 39755191

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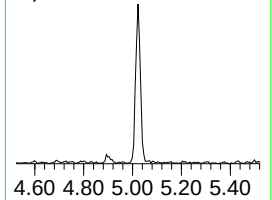
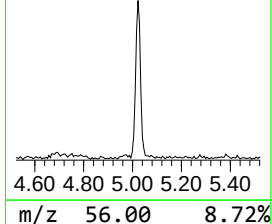
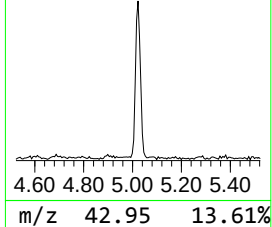
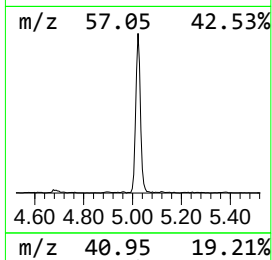
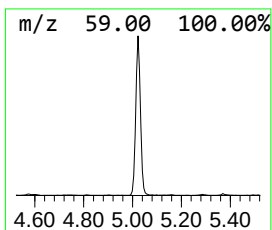
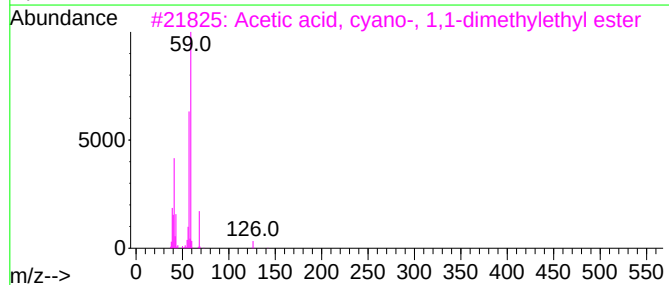
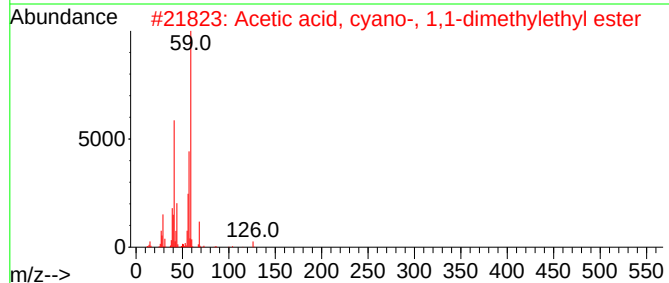
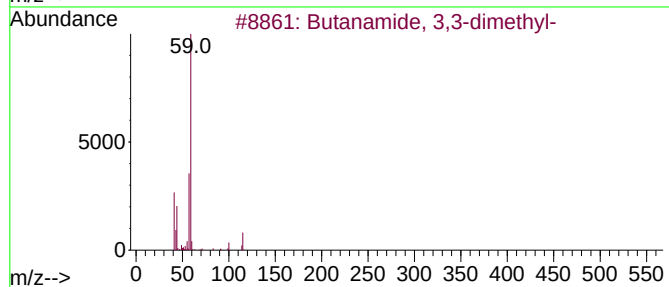
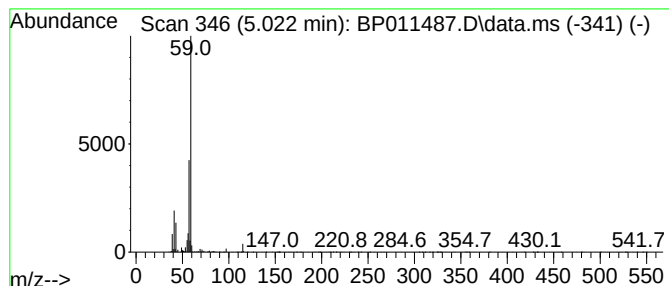
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TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Butanamide, 3,3-dimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.022	6.06 ng/ul	171826	1,4-Dichlorobenzene-d4	7.558

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butanamide, 3,3-dimethyl-	115	C6H13NO	000926-04-5	72
2			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	56
3			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	56
4			3-Methyl-6-ethyl--2,4-dioxadecane	188	C11H24O2	1000334-83-2	40
5			Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	40



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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
Butanamide, 3,3...	5.022	6.1	ng/ul	171826	1	7.558	567277	20.0