

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041923\
 Data File : BP014733.D
 Acq On : 19 Apr 2023 16:42
 Operator : CG/JU
 Sample : 02340-08
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BHE06

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-BP041823.M
 Title : SVOA CALIBRATION

Signal : TIC: BP014733.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.217	3	5	21	rVB	827214	895015	3.32%	1.176%
2	3.464	41	47	55	rBV	2620011	3710331	13.77%	4.874%
3	3.522	55	57	72	rVB	102694	152820	0.57%	0.201%
4	3.922	123	125	139	rBV	163722	261758	0.97%	0.344%
5	4.816	270	277	295	rBV	14952003	26940609	100.00%	35.391%
6	5.552	397	402	415	rVB	63426	97306	0.36%	0.128%
7	7.287	690	697	704	rBV	189921	292120	1.08%	0.384%
8	7.469	720	728	737	rBV	697365	1061088	3.94%	1.394%
9	7.675	755	763	774	rBV	644489	998550	3.71%	1.312%
10	8.152	837	844	857	rBV	542751	879317	3.26%	1.155%
11	8.852	953	963	970	rBV	503114	819349	3.04%	1.076%
12	9.322	1035	1043	1052	rBV	697510	1124704	4.17%	1.477%
13	10.051	1159	1167	1180	rBV	443658	720199	2.67%	0.946%
14	10.587	1250	1258	1272	rBV	861231	1426205	5.29%	1.874%
15	10.981	1316	1325	1334	rBV	757617	1257191	4.67%	1.652%
16	11.110	1340	1347	1356	rBV	522961	858804	3.19%	1.128%
17	12.551	1587	1592	1598	rVB2	18001	28272	0.10%	0.037%
18	14.187	1863	1870	1882	rBV	1792794	2549759	9.46%	3.350%
19	14.481	1913	1920	1927	rBV	2119735	3089000	11.47%	4.058%
20	14.787	1962	1972	1982	rBV	1150927	1718796	6.38%	2.258%
21	14.945	1993	1999	2005	rBV	132698	201560	0.75%	0.265%
22	15.775	2133	2140	2151	rBV	2758986	3920434	14.55%	5.150%
23	15.875	2151	2157	2168	rBV	423322	611521	2.27%	0.803%
24	16.134	2194	2201	2206	rBV	106385	152942	0.57%	0.201%
25	16.486	2256	2261	2269	rBV2	18036	31793	0.12%	0.042%
26	17.533	2433	2439	2448	rBV	1501202	2091356	7.76%	2.747%
27	17.633	2450	2456	2467	rVV2	3190561	4388148	16.29%	5.765%
28	18.392	2581	2585	2590	rBV	53391	74259	0.28%	0.098%
29	19.427	2757	2761	2766	rVB2	41545	53551	0.20%	0.070%
30	19.492	2768	2772	2776	rBV2	38327	56926	0.21%	0.075%
31	19.616	2790	2793	2797	rVB5	28228	32324	0.12%	0.042%
32	19.851	2827	2833	2841	rBV	4087262	5342535	19.83%	7.018%
33	21.286	3074	3077	3082	rBV3	72766	87731	0.33%	0.115%
34	21.610	3126	3132	3139	rBV	1679363	2280673	8.47%	2.996%
35	24.021	3534	3542	3551	rBV2	2719289	5529592	20.53%	7.264%
36	24.186	3563	3570	3581	rVB2	1153239	2386245	8.86%	3.135%

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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

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP041823.M
Title : SVOA CALIBRATION

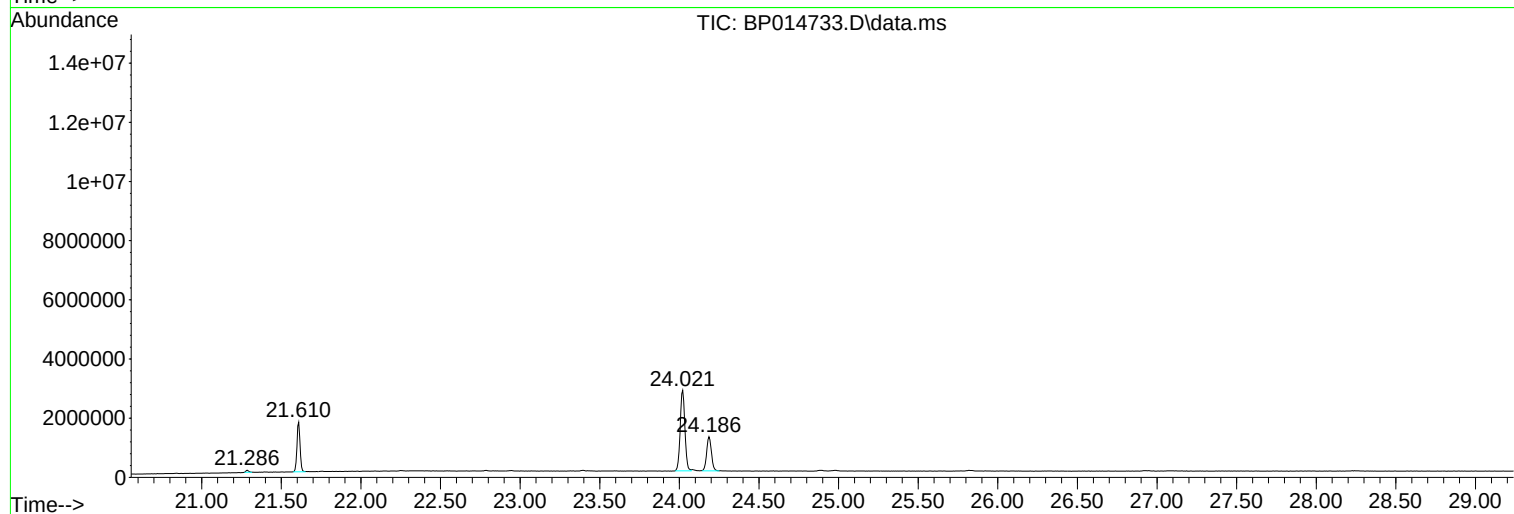
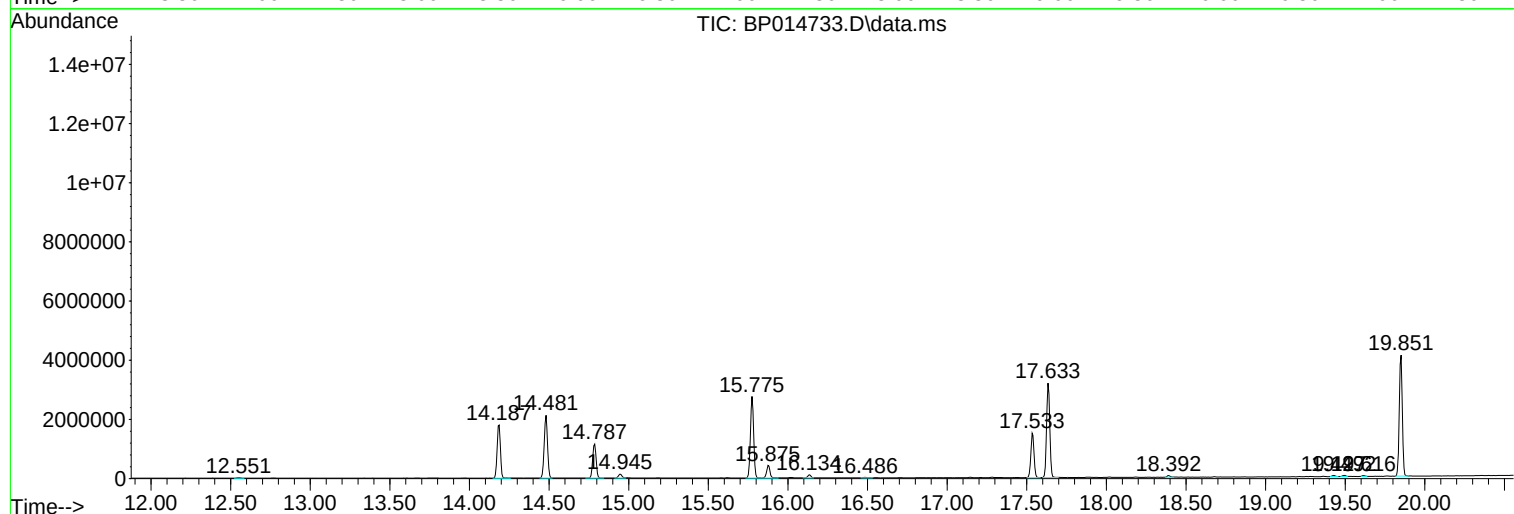
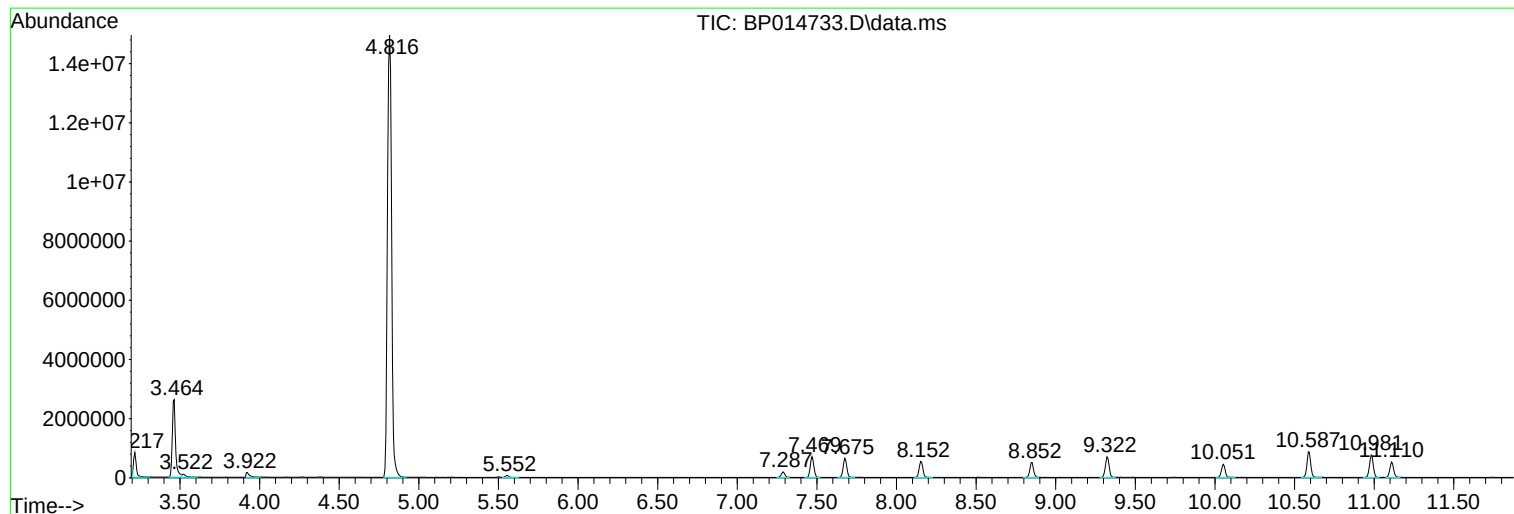
Sum of corrected areas: 76122783

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP041823.M
Quant Title : SVOA CALIBRATION

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



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Instrument :
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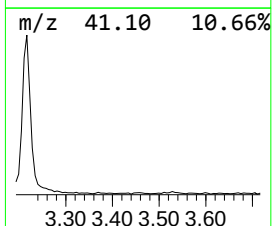
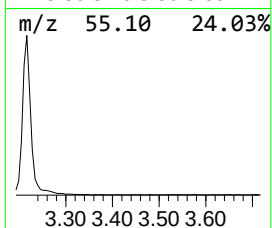
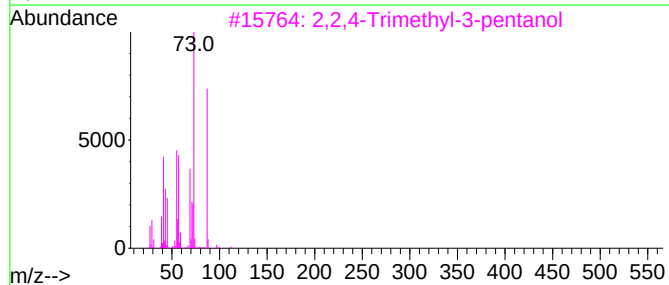
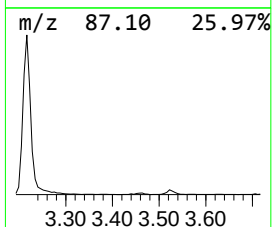
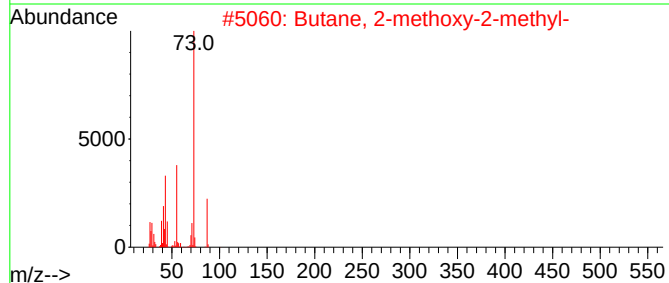
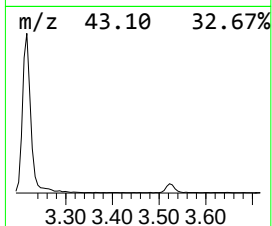
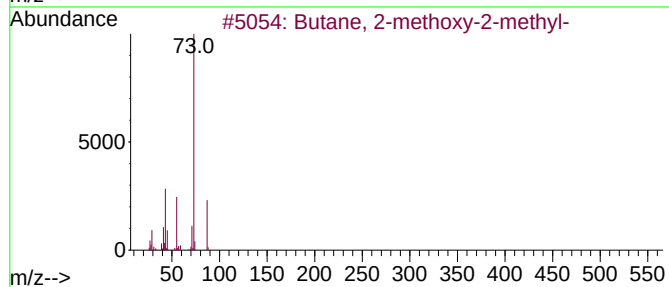
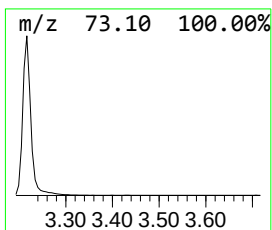
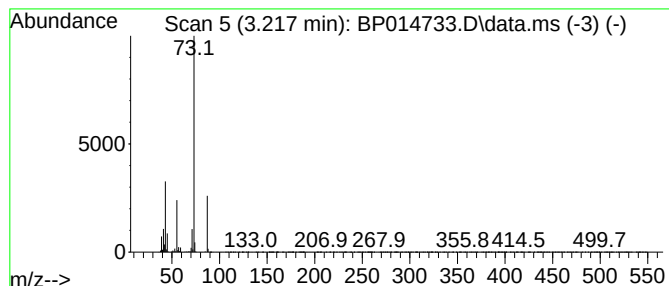
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP041823.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.217	20.36 ng/ul	895015	1,4-Dichlorobenzene-d4	8.152

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78		
2	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	74		
3	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39		
4	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23		
5	N-Ethylformamide	73	C3H7NO	000627-45-2	9		



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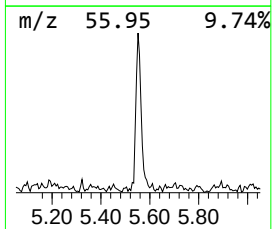
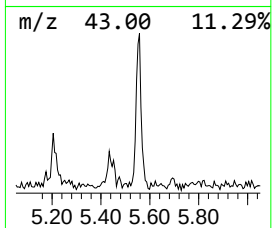
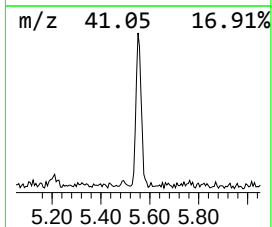
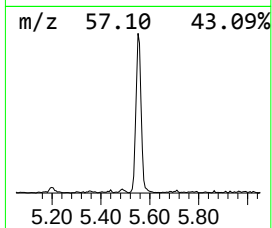
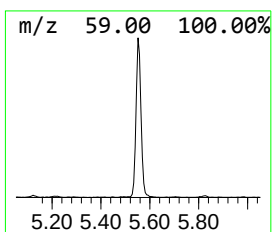
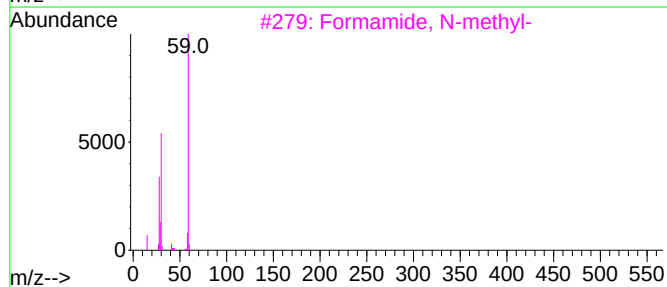
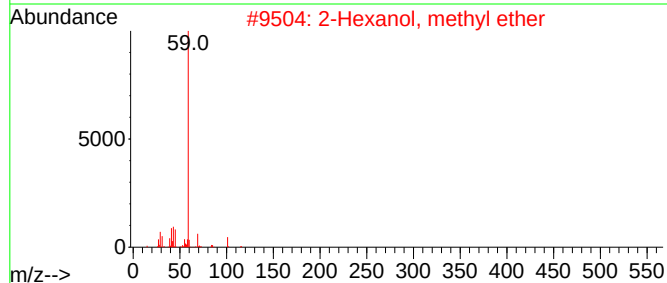
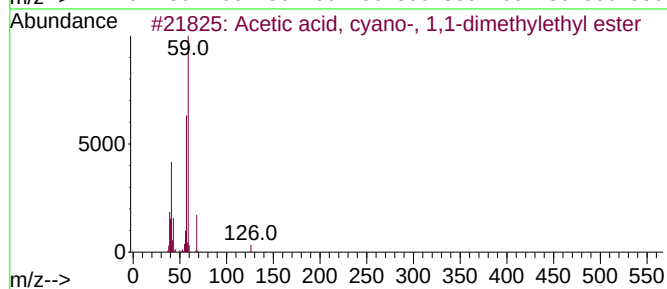
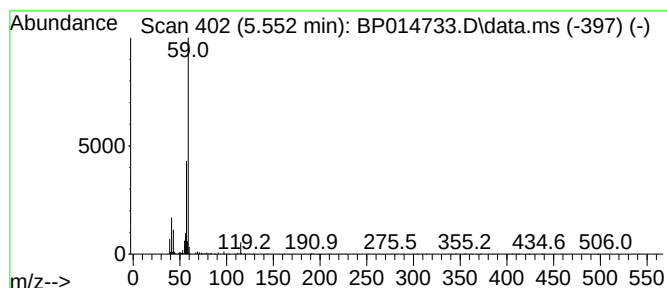
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP041823.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
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Peak Number 3 Acetic acid, cyano-, 1,1-di... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.552	2.21 ng/ul	97306	1,4-Dichlorobenzene-d4	8.152

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	56
2			2-Hexanol, methyl ether	116	C7H16O	1000333-92-0	45
3			Formamide, N-methyl-	59	C2H5NO	000123-39-7	9
4			Butanamide, 3,3-dimethyl-	115	C6H13NO	000926-04-5	9
5			3-Pentanol	88	C5H12O	000584-02-1	9



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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
Butane, 2-metho...	3.217	20.4	ng/ul	895015	1	8.152	879317	20.0
Acetic acid, cy...	5.552	2.2	ng/ul	97306	1	8.152	879317	20.0