

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG020322\
 Data File : BG052264.D
 Acq On : 3 Feb 2022 11:53
 Operator : CG/JU
 Sample : N1349-02
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0HP1

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_G\Methods\SFAM-EPA-BG013122.M
 Title : SVOA CALIBRATION

Signal : TIC: BG052264.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.540	3	9	36	rVB	11748813	24602393	100.00%	51.229%
2	3.998	83	87	96	rBV	24321	46474	0.19%	0.097%
3	4.345	140	146	155	rBV	46552	77312	0.31%	0.161%
4	4.439	157	162	169	rBV3	18884	31462	0.13%	0.066%
5	4.903	232	241	256	rBV	7969612	13845441	56.28%	28.830%
6	5.626	360	364	370	rBV	19623	29375	0.12%	0.061%
7	7.376	656	662	672	rVB2	38502	68870	0.28%	0.143%
8	7.541	684	690	698	rVB	117507	191842	0.78%	0.399%
9	7.764	720	728	736	rBV	116911	202440	0.82%	0.422%
10	8.234	800	808	815	rBV	112330	197368	0.80%	0.411%
11	8.939	921	928	937	rVB	99617	170168	0.69%	0.354%
12	9.403	999	1007	1017	rVB	148344	269548	1.10%	0.561%
13	10.137	1124	1132	1143	rBV	125872	220833	0.90%	0.460%
14	10.684	1219	1225	1235	rBV	165254	297256	1.21%	0.619%
15	11.071	1284	1291	1302	rVB	154823	274711	1.12%	0.572%
16	11.201	1305	1313	1322	rBV	148902	284839	1.16%	0.593%
17	14.267	1828	1835	1841	rBV	368597	511156	2.08%	1.064%
18	14.572	1880	1887	1898	rVB	388146	609184	2.48%	1.269%
19	14.878	1931	1939	1946	rBV	258426	389449	1.58%	0.811%
20	15.066	1965	1971	1978	rBV2	25839	49319	0.20%	0.103%
21	15.865	2101	2107	2115	rVB	510887	775508	3.15%	1.615%
22	15.976	2120	2126	2139	rVB	154928	227663	0.93%	0.474%
23	16.646	2233	2240	2247	rBV2	16336	25812	0.10%	0.054%
24	17.627	2400	2407	2413	rBV	318923	495188	2.01%	1.031%
25	17.727	2417	2424	2431	rVB	595380	888010	3.61%	1.849%
26	20.000	2805	2811	2819	rBV	738119	1060173	4.31%	2.208%
27	21.939	3134	3141	3150	rBV	321861	548608	2.23%	1.142%
28	25.176	3679	3692	3707	rVB2	351506	1031638	4.19%	2.148%
29	25.411	3721	3732	3744	rBV2	185933	572739	2.33%	1.193%
30	26.685	3940	3949	3953	rBV9	10566	29151	0.12%	0.061%

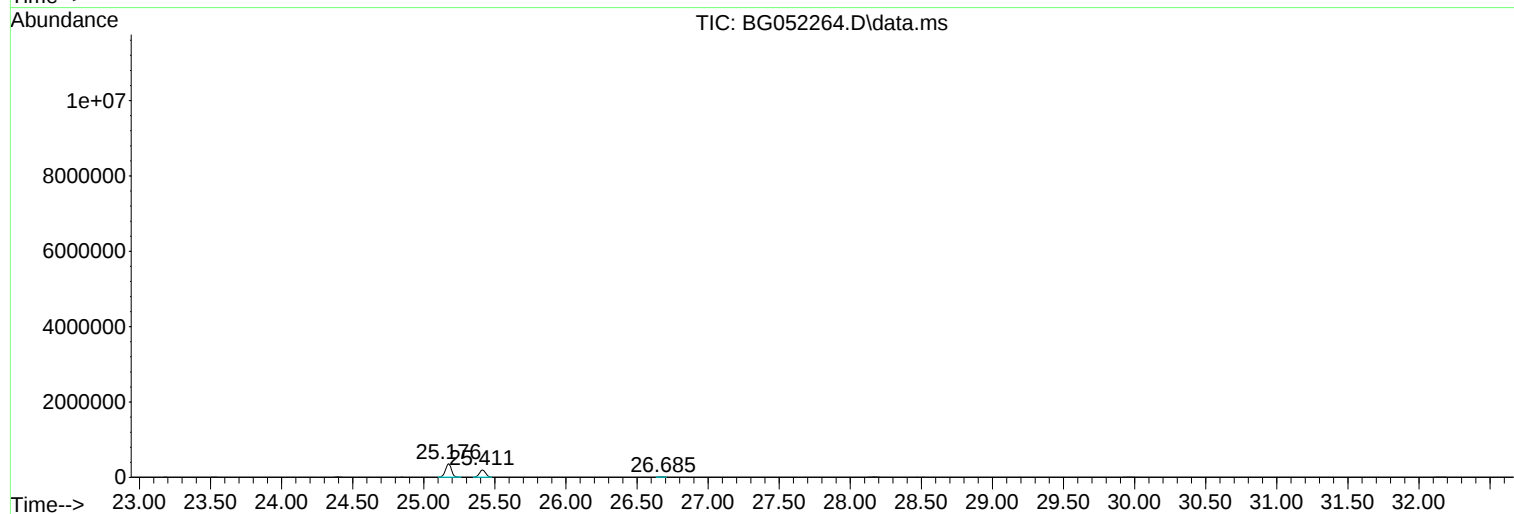
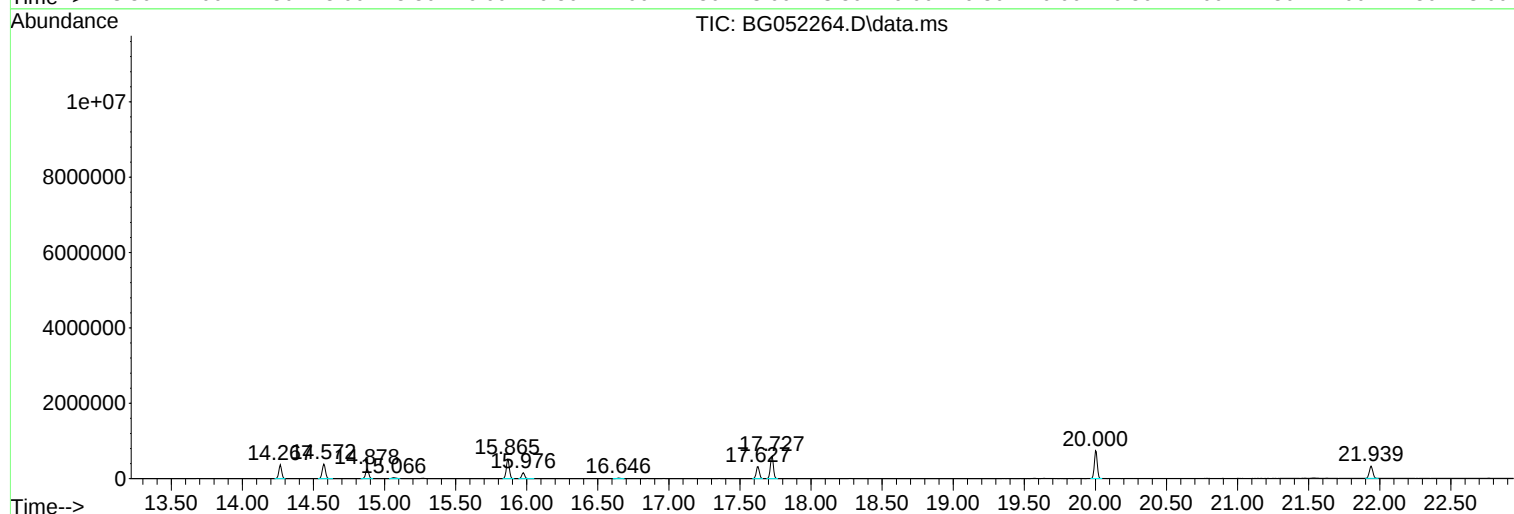
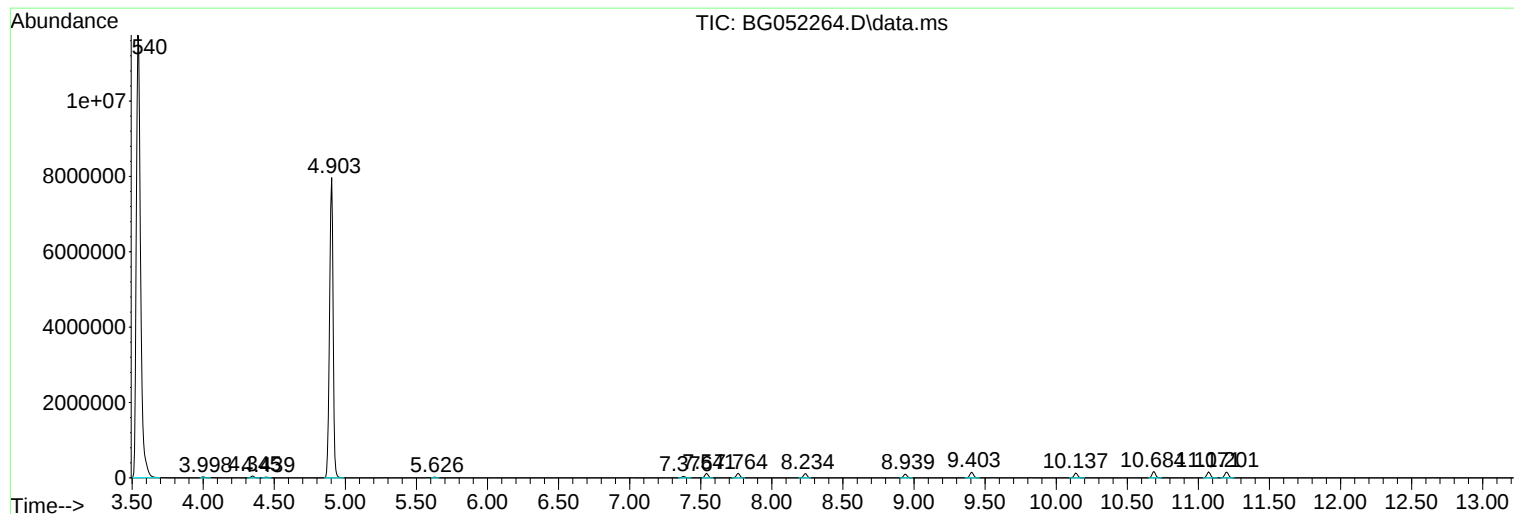
Sum of corrected areas: 48023930

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG013122.M
Quant Title : SVOA CALIBRATION

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



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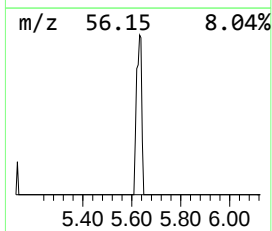
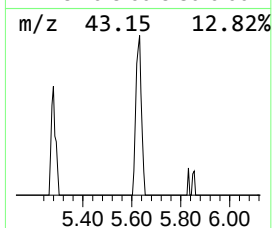
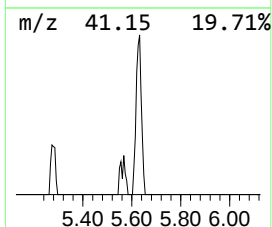
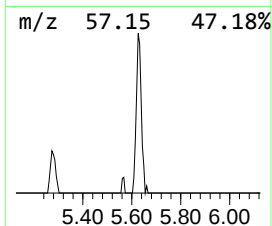
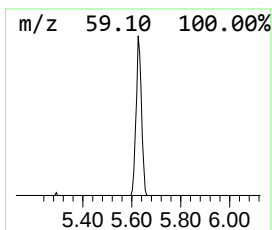
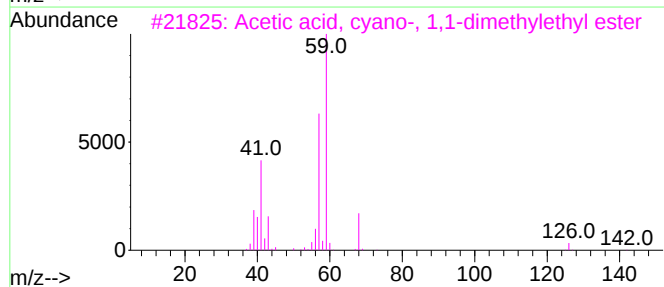
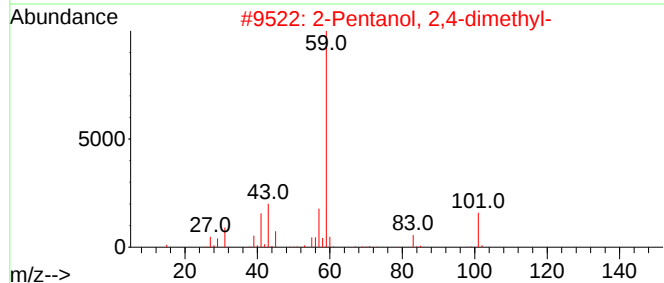
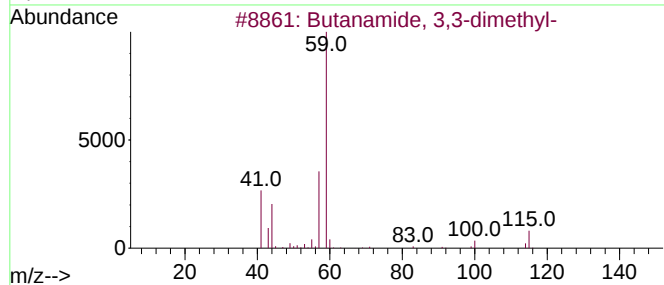
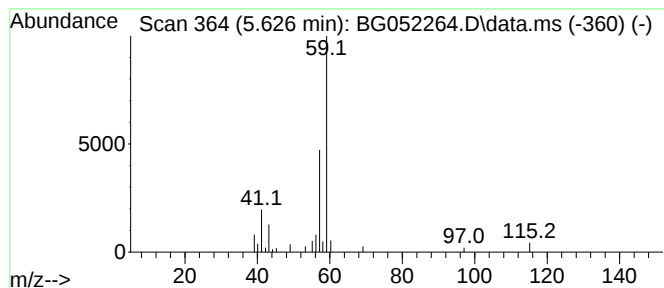
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG013122.M
Quant Title : SVOA CALIBRATION

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

Peak Number 4 Butanamide, 3,3-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.626	2.98 ng/ul	29375	1,4-Dichlorobenzene-d4	8.234

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butanamide, 3,3-dimethyl-	115	C6H13NO	000926-04-5	72
2			2-Pentanol, 2,4-dimethyl-	116	C7H16O	000625-06-9	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\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butanamide, 3,3...	5.626	3.0	ng/ul	29375	1	8.234	197368	20.0