

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG012622\
 Data File : BG052195.D
 Acq On : 27 Jan 2022 21:44
 Operator : CG/JU
 Sample : N1273-04
 Misc :
 ALS Vial : 54 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0HK3

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-BG010622.M
 Title : SVOA CALIBRATION

Signal : TIC: BG052195.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.541	3	9	16	rBV	981282	1763059	19.12%	8.540%
2	3.599	16	19	34	rVB	327493	523267	5.67%	2.534%
3	4.005	81	88	92	rBV4	10701	19882	0.22%	0.096%
4	4.451	158	164	173	rBV5	4786	9926	0.11%	0.048%
5	4.903	232	241	253	rBV	5500823	9223169	100.00%	44.673%
6	4.992	253	256	261	rVB3	7888	10461	0.11%	0.051%
7	7.382	653	663	673	rVB3	25244	50412	0.55%	0.244%
8	7.541	683	690	698	rBV	91805	152155	1.65%	0.737%
9	7.764	722	728	738	rBV	95008	165738	1.80%	0.803%
10	7.999	762	768	774	rBV2	13098	19187	0.21%	0.093%
11	8.240	798	809	815	rBV	93293	159053	1.72%	0.770%
12	8.939	921	928	935	rBV2	73041	126984	1.38%	0.615%
13	9.409	999	1008	1017	rBV	120869	219771	2.38%	1.064%
14	9.826	1075	1079	1084	rVB2	7666	11323	0.12%	0.055%
15	10.138	1124	1132	1140	rVB	103102	182023	1.97%	0.882%
16	10.690	1218	1226	1233	rBV	139594	257534	2.79%	1.247%
17	11.077	1284	1292	1299	rBV	134161	235633	2.55%	1.141%
18	11.201	1306	1313	1322	rBV	99797	180463	1.96%	0.874%
19	14.267	1829	1835	1841	rBV	314066	467309	5.07%	2.263%
20	14.496	1865	1874	1881	rBV	383051	728857	7.90%	3.530%
21	14.573	1881	1887	1897	rVB	331616	542703	5.88%	2.629%
22	14.878	1933	1939	1948	rVB	227861	356761	3.87%	1.728%
23	15.066	1966	1971	1981	rBV3	22915	44179	0.48%	0.214%
24	15.871	2099	2108	2116	rBV	453843	681478	7.39%	3.301%
25	15.977	2120	2126	2137	rVB	125383	186010	2.02%	0.901%
26	17.627	2400	2407	2415	rBV	313816	466132	5.05%	2.258%
27	17.727	2417	2424	2434	rVB	512568	780056	8.46%	3.778%
28	18.279	2513	2518	2523	rBV4	9620	12859	0.14%	0.062%
29	19.642	2746	2750	2753	rBV2	6538	9241	0.10%	0.045%
30	19.889	2789	2792	2799	rBV5	7825	11106	0.12%	0.054%
31	20.007	2804	2812	2824	rVV	670347	937715	10.17%	4.542%
32	21.546	3070	3074	3077	rBV6	5921	10621	0.12%	0.051%
33	21.945	3135	3142	3151	rBV	301649	515134	5.59%	2.495%
34	23.190	3349	3354	3361	rVB8	11478	20045	0.22%	0.097%
35	23.537	3409	3413	3418	rBV6	7938	11055	0.12%	0.054%
36	24.412	3558	3562	3569	rVB6	9722	21405	0.23%	0.104%

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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	25.188	3682	3694	3708	rVB3	322805	918492	9.96%	4.449%
38	25.429	3723	3735	3748	rBV2	175457	548576	5.95%	2.657%
39	26.692	3944	3950	3961	rVB6	14243	41261	0.45%	0.200%
40	28.195	4197	4206	4207	rBV7	11384	24852	0.27%	0.120%

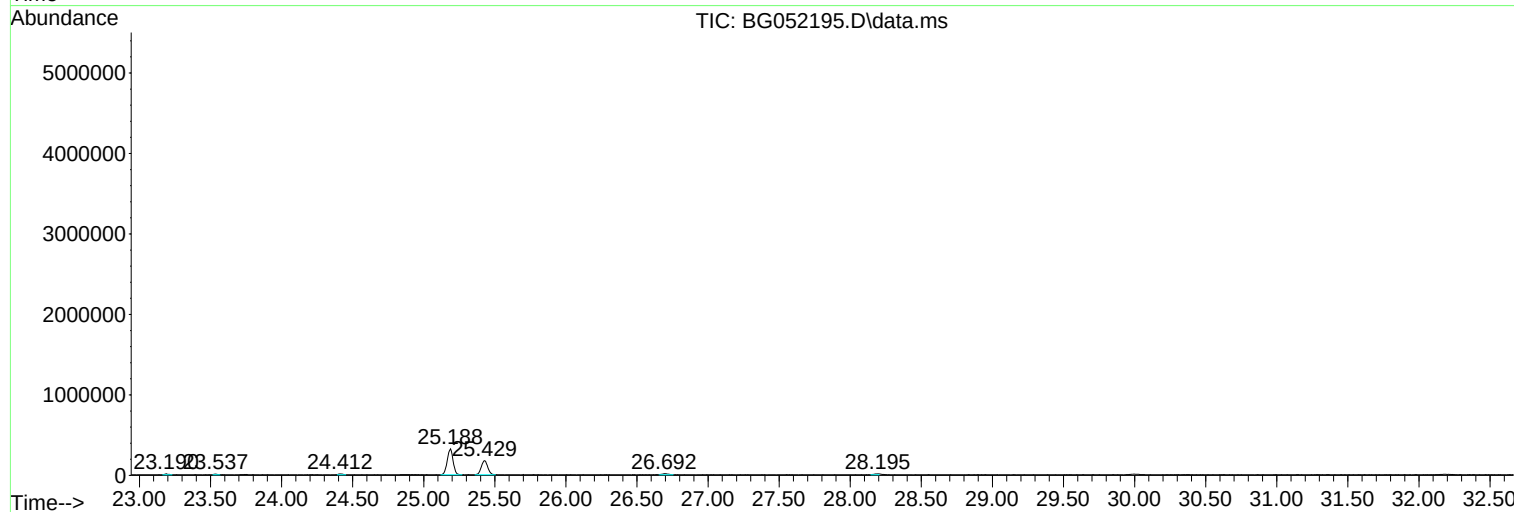
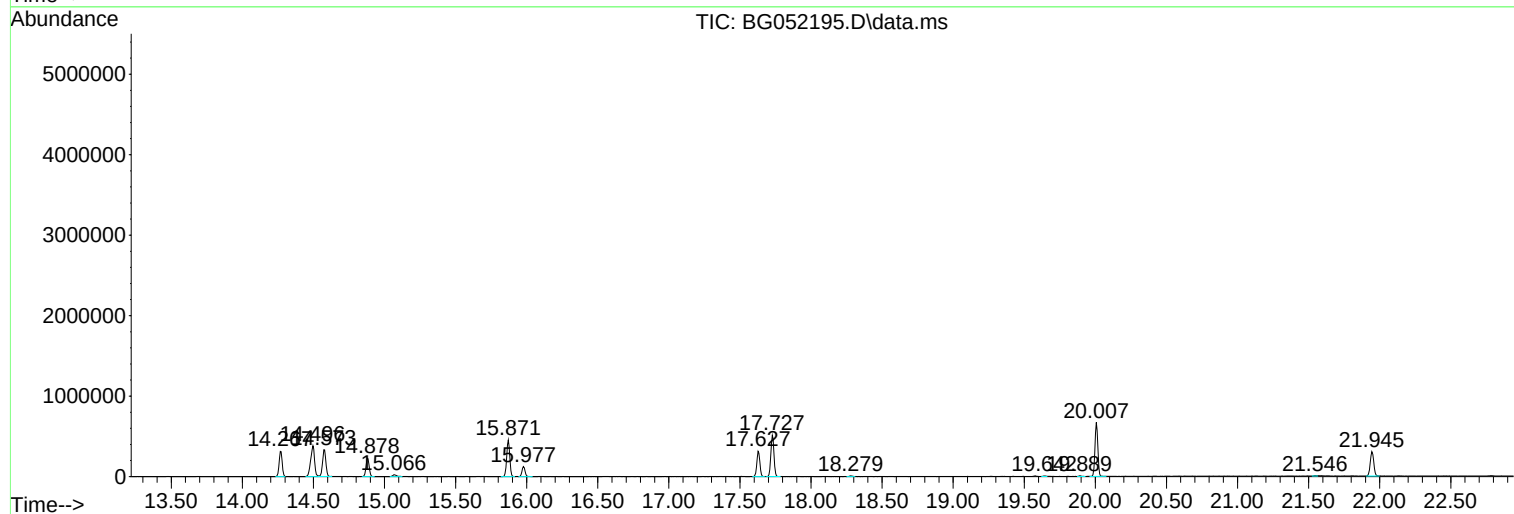
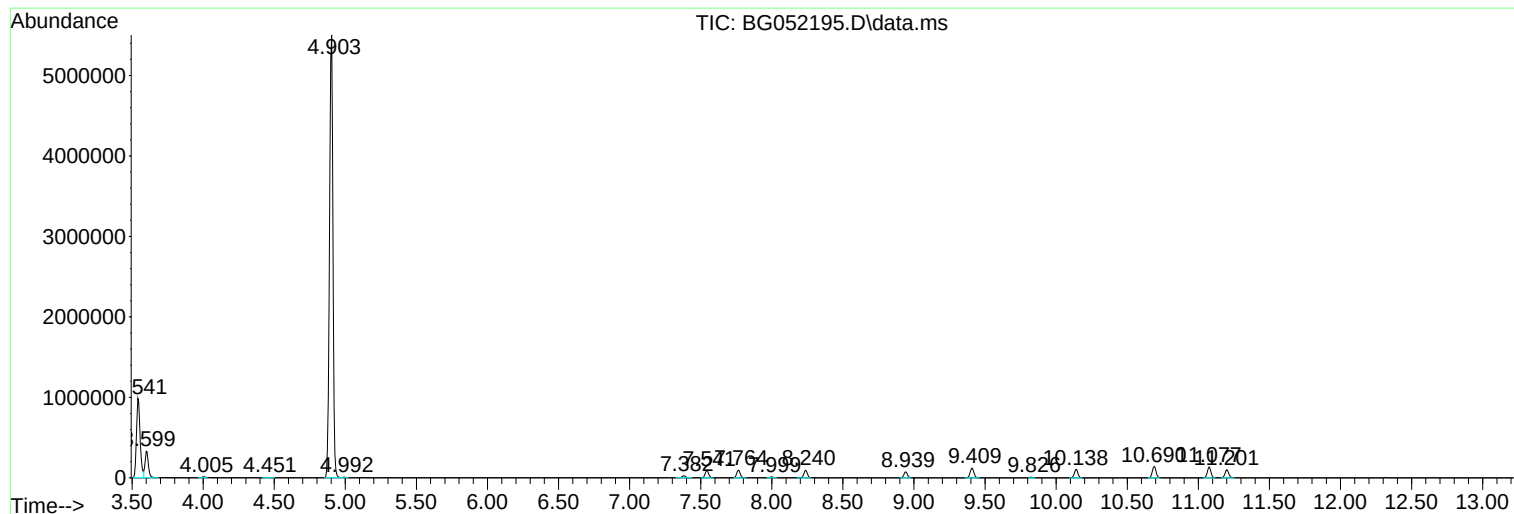
Sum of corrected areas: 20645887

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG010622.M
Quant Title : SVOA CALIBRATION

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



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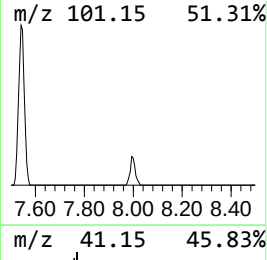
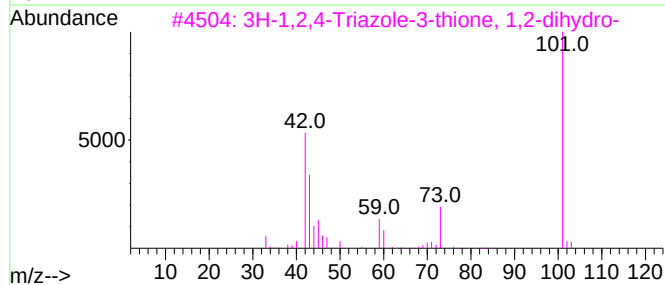
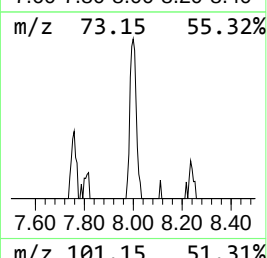
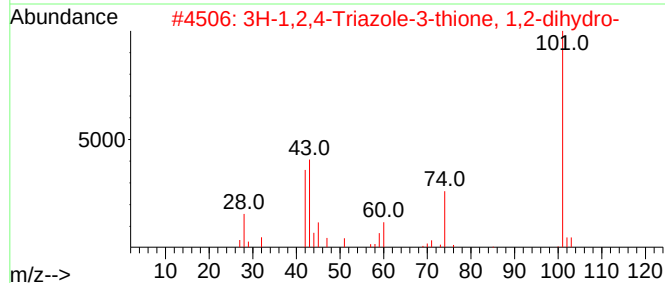
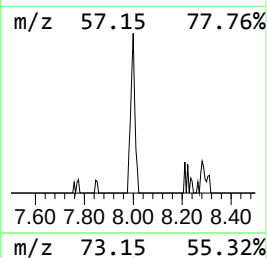
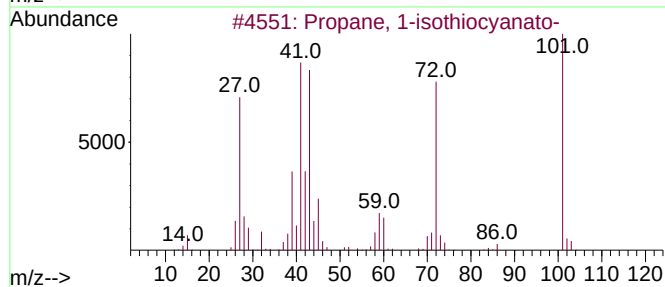
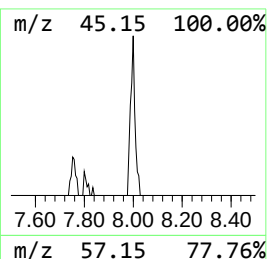
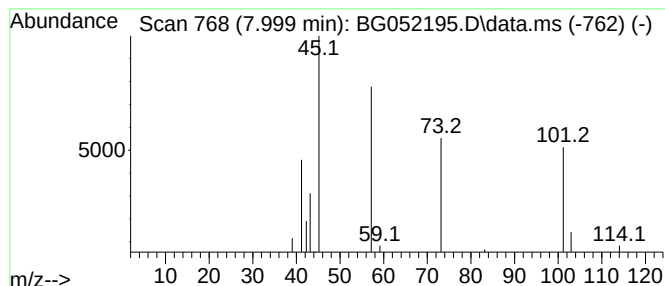
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Peak Number 2 unknown-01 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.999	2.41 ng/ul	19187	1,4-Dichlorobenzene-d4	8.240

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propane, 1-isothiocyanato-	101	C4H7NS	000628-30-8	12
2			3H-1,2,4-Triazole-3-thione, 1,2-...	101	C2H3N3S	003179-31-5	9
3			3H-1,2,4-Triazole-3-thione, 1,2-...	101	C2H3N3S	003179-31-5	9
4			2-Pyrrolidinethione	101	C4H7NS	002295-35-4	9
5			Propane, 1-isothiocyanato-	101	C4H7NS	000628-30-8	9



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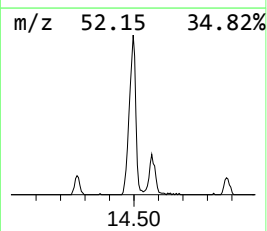
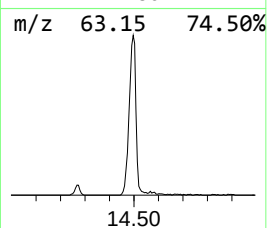
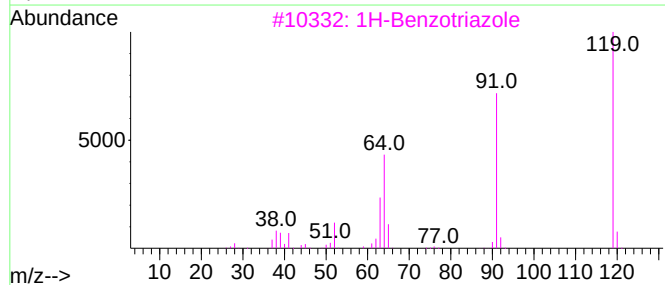
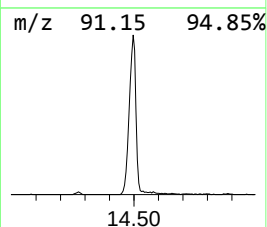
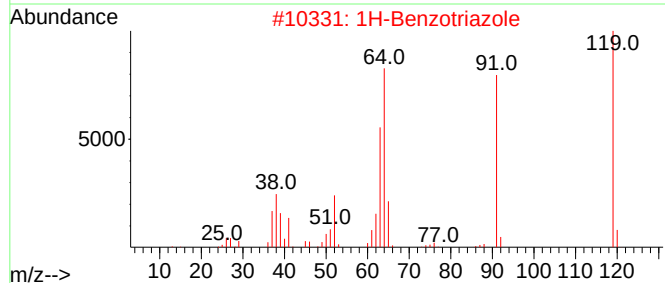
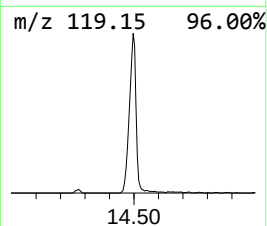
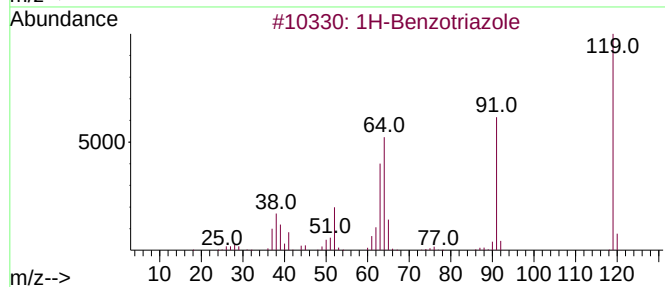
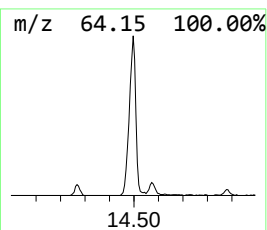
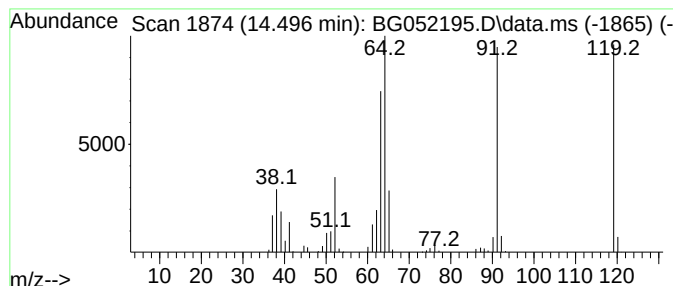
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Peak Number 3 1H-Benzotriazole Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.496	40.86 ng/ul	728857	Acenaphthene-d10	14.884

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Benzotriazole			119	C6H5N3	000095-14-7	98
2	1H-Benzotriazole			119	C6H5N3	000095-14-7	98
3	1H-Benzotriazole			119	C6H5N3	000095-14-7	96
4	1H-Benzotriazole			119	C6H5N3	000095-14-7	95
5	Ethane-1,2-diol, 1,2-bis(1-benzo...			296	C14H12N6O2	111507-88-1	91



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TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown-01	7.999	2.4	ng/ul	19187	1	8.240	159053	20.0
1H-Benzotriazole	14.496	40.9	ng/ul	728857	3	14.884	356761	20.0