

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG070824\
 Data File : BG061939.D
 Acq On : 8 Jul 2024 12:44
 Operator : MA/JU
 Sample : P3059-02
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 E1GM4

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-BG070224.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BG061939.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.463	26	30	31	rBV4	5885	6886	0.31%	0.033%
2	3.480	31	33	37	rVB2	11102	11409	0.51%	0.055%
3	3.627	56	58	60	rVB3	3377	3222	0.14%	0.016%
4	3.757	77	80	87	rVB3	3935	7012	0.31%	0.034%
5	3.904	102	105	112	rBV	49169	81951	3.65%	0.398%
6	3.956	112	114	117	rVB3	4478	4689	0.21%	0.023%
7	5.161	315	319	320	rBV2	4419	6697	0.30%	0.032%
8	5.331	344	348	350	rVB2	3002	3051	0.14%	0.015%
9	6.512	545	549	551	rBV2	3191	2886	0.13%	0.014%
10	6.700	579	581	585	rBV	2679	3372	0.15%	0.016%
11	6.771	592	593	599	rVB	2101	2857	0.13%	0.014%
12	6.835	599	604	605	rBV2	2817	3606	0.16%	0.017%
13	6.882	611	612	618	rVB2	2313	3360	0.15%	0.016%
14	7.247	668	674	680	rBV2	85277	149982	6.68%	0.728%
15	7.417	696	703	710	rVB2	187071	309993	13.81%	1.504%
16	7.623	731	738	745	rVV2	244289	407170	18.14%	1.976%
17	7.693	748	750	755	rVB3	4047	5694	0.25%	0.028%
18	7.869	777	780	784	rBV	1936	2991	0.13%	0.015%
19	7.993	797	801	804	rBV	2219	3657	0.16%	0.018%
20	8.093	811	818	824	rBV	185670	321106	14.31%	1.558%
21	8.140	824	826	831	rVV3	3634	3856	0.17%	0.019%
22	8.222	835	840	841	rBV2	2401	3990	0.18%	0.019%
23	8.604	902	905	909	rBV2	2993	3851	0.17%	0.019%
24	8.798	929	938	946	rVB	249518	430453	19.18%	2.089%
25	8.880	951	952	956	rVB	3283	3564	0.16%	0.017%
26	8.962	963	966	968	rBV	2795	3038	0.14%	0.015%
27	9.168	997	1001	1003	rBV	2281	3895	0.17%	0.019%
28	9.262	1008	1017	1024	rVV	307094	558639	24.89%	2.711%
29	9.503	1055	1058	1060	rBV	2902	2998	0.13%	0.015%
30	9.985	1133	1140	1147	rBV	279620	505560	22.53%	2.453%
31	10.525	1223	1232	1242	rVB	381313	700288	31.21%	3.398%
32	10.907	1287	1297	1305	rBV	299591	539355	24.03%	2.617%
33	11.042	1310	1320	1327	rBV	314322	575252	25.63%	2.792%
34	11.377	1372	1377	1380	rBV	1894	3488	0.16%	0.017%
35	11.577	1408	1411	1414	rVB	3212	3241	0.14%	0.016%
36	11.636	1417	1421	1427	rVV3	9449	18408	0.82%	0.089%

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Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
 Title : SVOA CALIBRATION

37	11.694	1429	1431	1436	rVB	3372	3316	0.15%	0.016%
38	12.388	1543	1549	1551	rBV3	5918	8307	0.37%	0.040%
39	12.476	1560	1564	1570	rBV4	6341	11794	0.53%	0.057%
40	12.623	1586	1589	1591	rBV2	2448	2937	0.13%	0.014%
41	12.987	1649	1651	1655	rVB	3546	3833	0.17%	0.019%
42	13.140	1674	1677	1680	rVB	2624	3307	0.15%	0.016%
43	13.228	1688	1692	1695	rBV2	2566	3363	0.15%	0.016%
44	13.892	1801	1805	1807	rBV	2514	3396	0.15%	0.016%
45	14.115	1837	1843	1849	rBV	843655	1223117	54.50%	5.936%
46	14.268	1867	1869	1872	rBV2	2740	3320	0.15%	0.016%
47	14.415	1887	1894	1902	rVB	839732	1310462	58.40%	6.359%
48	14.509	1907	1910	1912	rBV3	3511	2902	0.13%	0.014%
49	14.720	1939	1946	1953	rVB	522994	837890	37.34%	4.066%
50	14.902	1972	1977	1986	rBV2	106552	186972	8.33%	0.907%
51	15.120	2011	2014	2016	rBV2	3448	3071	0.14%	0.015%
52	15.226	2030	2032	2035	rBV2	3002	2976	0.13%	0.014%
53	15.525	2081	2083	2086	rVB	3145	3036	0.14%	0.015%
54	15.602	2094	2096	2099	rBV	3257	3847	0.17%	0.019%
55	15.707	2107	2114	2122	rVB	1082970	1668803	74.36%	8.098%
56	15.819	2125	2133	2139	rBV	514130	727043	32.40%	3.528%
57	15.948	2151	2155	2158	rBV2	7403	12731	0.57%	0.062%
58	16.036	2167	2170	2172	rBV2	2821	3442	0.15%	0.017%
59	16.095	2179	2180	2185	rVB	3657	3843	0.17%	0.019%
60	16.700	2281	2283	2285	rBV2	3335	3454	0.15%	0.017%
61	17.088	2346	2349	2354	rBV2	2643	4288	0.19%	0.021%
62	17.200	2366	2368	2371	rVB2	3671	2960	0.13%	0.014%
63	17.235	2371	2374	2377	rBV3	3186	3721	0.17%	0.018%
64	17.364	2393	2396	2399	rVB3	4602	5499	0.25%	0.027%
65	17.458	2405	2412	2418	rBV	716745	1057652	47.13%	5.133%
66	17.558	2423	2429	2436	rVB	1293549	1910474	85.13%	9.271%
67	17.899	2485	2487	2492	rBV3	3536	4245	0.19%	0.021%
68	18.069	2515	2516	2519	rVB	3914	3643	0.16%	0.018%
69	18.099	2519	2521	2523	rBV	3772	4215	0.19%	0.020%
70	18.363	2564	2566	2570	rBV3	3597	4562	0.20%	0.022%
71	18.404	2570	2573	2579	rVB4	3970	7836	0.35%	0.038%
72	18.557	2597	2599	2603	rBV2	3307	4581	0.20%	0.022%
73	18.710	2623	2625	2628	rVB3	3901	3095	0.14%	0.015%
74	18.804	2638	2641	2645	rBV4	4664	7648	0.34%	0.037%
75	18.898	2655	2657	2659	rBV3	4996	4292	0.19%	0.021%
76	18.974	2669	2670	2674	rBV3	5522	5578	0.25%	0.027%

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Title : SVOA CALIBRATION

77	19.397	2738	2742	2747	rVB4	18448	27709	1.23%	0.134%
78	19.468	2751	2754	2759	rVB3	17936	26616	1.19%	0.129%
79	19.744	2799	2801	2805	rVB3	10495	10522	0.47%	0.051%
80	19.838	2810	2817	2824	rVB	1462837	2080297	92.70%	10.095%
81	20.402	2911	2913	2914	rBV2	8518	4048	0.18%	0.020%
82	21.001	3013	3015	3016	rBV2	8719	4746	0.21%	0.023%
83	21.718	3130	3137	3144	rBV	638861	1081988	48.21%	5.251%
84	22.517	3269	3273	3282	rBV	18713	40538	1.81%	0.197%
85	22.893	3335	3337	3339	rVB2	306630	112395	5.01%	0.545%
86	24.738	3642	3651	3665	rVB2	788061	2244121	100.00%	10.890%
87	24.961	3680	3689	3700	rVB2	407302	1200709	53.50%	5.827%

Sum of corrected areas: 20606607

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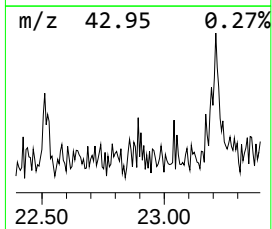
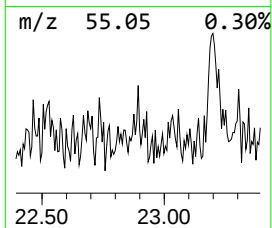
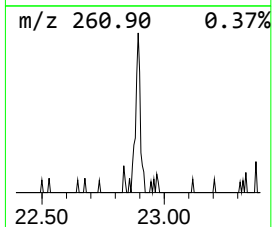
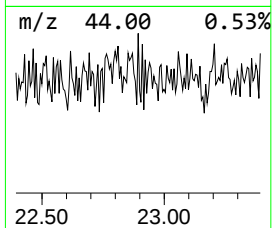
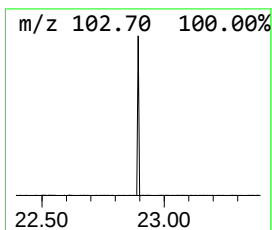
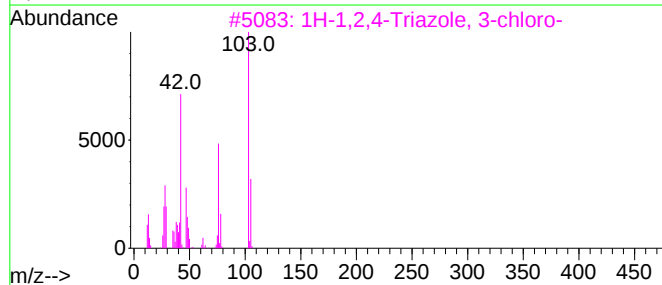
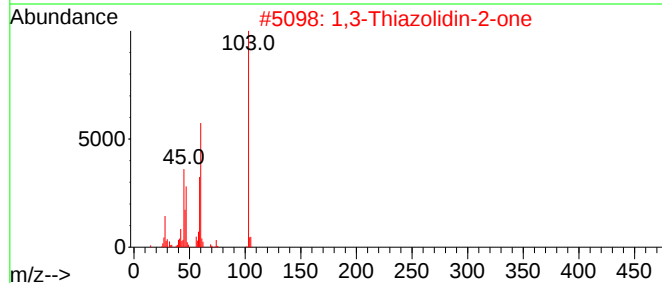
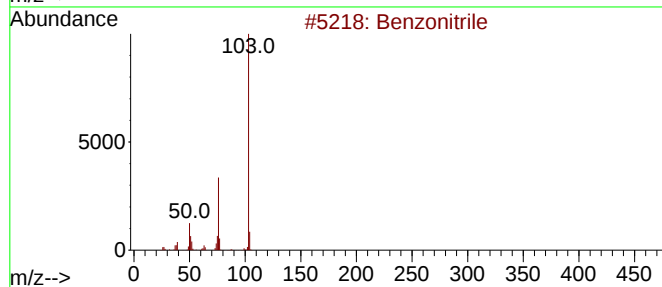
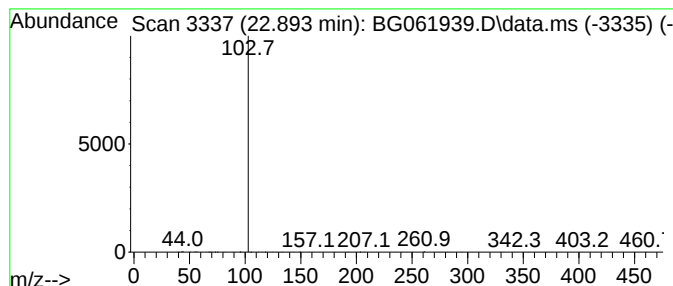
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 unknown-01 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.893	2.08 ng/ul	112395	Chrysene-d12	21.718

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzonitrile	103	C7H5N	000100-47-0	2
2			1,3-Thiazolidin-2-one	103	C3H5NOS	002682-49-7	2
3			1H-1,2,4-Triazole, 3-chloro-	103	C2H2ClN3	006818-99-1	2
4			Ethanethioamide, N,N-dimethyl-	103	C4H9NS	000631-67-4	2
5			Ethanethioamide, N,N-dimethyl-	103	C4H9NS	000631-67-4	2



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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
unknown-01	22.893	2.1	ng/u1	112395	5	21.718	1081990	20.0