

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071823\
 Data File : BG058354.D
 Acq On : 20 Jul 2023 9:59
 Operator : CG/JU
 Sample : 03452-18
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 A46Q6

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-BG071723.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BG058354.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.152	3	11	36	rVB	1927388	3310774	100.00%	24.312%
2	3.340	40	43	46	rVB4	4878	6748	0.20%	0.050%
3	3.751	108	113	125	rBV	77290	130834	3.95%	0.961%
4	4.086	165	170	176	rVB2	15308	23843	0.72%	0.175%
5	4.150	177	181	191	rBV	37426	74303	2.24%	0.546%
6	4.327	209	211	216	rVB4	5689	7159	0.22%	0.053%
7	4.609	257	259	266	rBV5	2797	3661	0.11%	0.027%
8	5.355	381	386	391	rVB4	3494	5670	0.17%	0.042%
9	5.796	453	461	466	rBV	115200	212134	6.41%	1.558%
10	5.937	483	485	490	rVB5	3036	3867	0.12%	0.028%
11	6.001	493	496	502	rBV	2933	3791	0.11%	0.028%
12	6.172	521	525	532	rBV2	15565	25406	0.77%	0.187%
13	6.518	577	584	593	rBV	193053	322500	9.74%	2.368%
14	7.059	669	676	689	rBV	90469	166114	5.02%	1.220%
15	7.223	697	704	714	rBV	75787	126518	3.82%	0.929%
16	7.429	732	739	747	rBV	94230	164558	4.97%	1.208%
17	7.640	773	775	782	rVB6	2561	4251	0.13%	0.031%
18	7.893	811	818	825	rBV	128473	223099	6.74%	1.638%
19	7.969	827	831	837	rVB4	6338	11320	0.34%	0.083%
20	8.069	842	848	855	rBV4	10513	20453	0.62%	0.150%
21	8.146	855	861	866	rBV2	15837	30445	0.92%	0.224%
22	8.210	866	872	880	rVB	163841	279535	8.44%	2.053%
23	8.598	932	938	947	rBV	105452	198056	5.98%	1.454%
24	8.657	947	948	954	rVB3	4994	7075	0.21%	0.052%
25	8.892	983	988	993	rBV5	2812	5275	0.16%	0.039%
26	9.056	1010	1016	1027	rBV	94711	169641	5.12%	1.246%
27	9.186	1035	1038	1046	rVB5	3382	6788	0.21%	0.050%
28	9.779	1133	1139	1149	rVB2	78488	143736	4.34%	1.055%
29	9.944	1163	1167	1171	rBV6	3078	4736	0.14%	0.035%
30	10.320	1224	1231	1245	rBV	129236	248199	7.50%	1.823%
31	10.696	1286	1295	1303	rBV	210816	371944	11.23%	2.731%
32	10.837	1313	1319	1328	rBV	35009	72640	2.19%	0.533%
33	11.160	1368	1374	1382	rVB3	2717	4819	0.15%	0.035%
34	11.935	1502	1506	1514	rVB4	2087	3887	0.12%	0.029%
35	12.282	1561	1565	1570	rVB2	3371	5310	0.16%	0.039%
36	12.746	1639	1644	1647	rBV2	8075	12588	0.38%	0.092%

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Filtering: 5

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Max Peaks: 100

Stop Thrs : 0

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If leading or trailing edge < 100 prefer < Baseline drop else tangent >

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37	12.781	1647	1650	1656	rVB2	6953	10835	0.33%	0.080%
38	12.893	1663	1669	1674	rBV4	2384	4164	0.13%	0.031%
39	13.140	1706	1711	1716	rVB5	2967	5222	0.16%	0.038%
40	13.422	1754	1759	1765	rBV3	3893	6546	0.20%	0.048%
41	13.933	1840	1846	1852	rBV	259458	375980	11.36%	2.761%
42	14.227	1890	1896	1909	rBV	294769	464483	14.03%	3.411%
43	14.391	1918	1924	1928	rVB3	2273	3576	0.11%	0.026%
44	14.532	1941	1948	1959	rBV	365147	566543	17.11%	4.160%
45	14.744	1977	1984	1997	rBV3	27104	68573	2.07%	0.504%
46	14.891	2003	2009	2015	rBV6	7228	15054	0.45%	0.111%
47	15.314	2075	2081	2084	rBV5	2301	4105	0.12%	0.030%
48	15.343	2084	2086	2089	rVB3	3570	3355	0.10%	0.025%
49	15.525	2109	2117	2125	rBV	411822	619792	18.72%	4.551%
50	15.643	2131	2137	2146	rVB2	44488	71016	2.14%	0.521%
51	15.766	2152	2158	2161	rVB4	2688	4287	0.13%	0.031%
52	15.884	2174	2178	2185	rVB	26891	37666	1.14%	0.277%
53	16.965	2356	2362	2370	rBV	93821	130636	3.95%	0.959%
54	17.153	2390	2394	2397	rBV	3367	4463	0.13%	0.033%
55	17.276	2406	2415	2426	rBV	470139	707556	21.37%	5.196%
56	17.376	2426	2432	2445	rVV2	394550	584655	17.66%	4.293%
57	18.158	2561	2565	2574	rBV2	34203	59819	1.81%	0.439%
58	18.234	2574	2578	2580	rVV	4964	7892	0.24%	0.058%
59	18.263	2580	2583	2589	rVB	20622	26525	0.80%	0.195%
60	18.663	2647	2651	2655	rBV6	2902	4674	0.14%	0.034%
61	18.821	2674	2678	2683	rVB3	8208	11993	0.36%	0.088%
62	19.233	2743	2748	2751	rBV4	6299	9452	0.29%	0.069%
63	19.297	2755	2759	2763	rBV2	10850	19881	0.60%	0.146%
64	19.333	2763	2765	2769	rVB2	9143	10686	0.32%	0.078%
65	19.432	2779	2782	2786	rBV4	7677	11973	0.36%	0.088%
66	19.579	2804	2807	2812	rVB6	3795	5603	0.17%	0.041%
67	19.668	2816	2822	2831	rVV	526319	775098	23.41%	5.692%
68	19.838	2849	2851	2856	rVV6	3287	4095	0.12%	0.030%
69	20.055	2885	2888	2890	rBV4	4363	5759	0.17%	0.042%
70	20.190	2907	2911	2915	rBV5	7857	9633	0.29%	0.071%
71	20.684	2992	2995	3000	rVB3	9496	9605	0.29%	0.071%
72	20.995	3046	3048	3049	rBV2	5954	3711	0.11%	0.027%
73	21.113	3065	3068	3072	rVB6	8347	10903	0.33%	0.080%
74	21.172	3075	3078	3081	rBV2	22435	29391	0.89%	0.216%
75	21.530	3132	3139	3154	rVB	415744	714190	21.57%	5.244%
76	21.700	3164	3168	3173	rBV2	18818	30281	0.91%	0.222%

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Sampling : 1 Min Area: 0.1 % of largest Peak
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Stop Thrs : 0 Peak Location: TOP

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77	22.288	3264	3268	3274	rVB4	24630	44027	1.33%	0.323%
78	22.964	3378	3383	3389	rVB2	20098	35344	1.07%	0.260%
79	23.733	3509	3514	3521	rVB2	27872	60557	1.83%	0.445%
80	24.444	3625	3635	3649	rBV2	249596	699100	21.12%	5.134%
81	24.662	3662	3672	3685	rVB	288432	810362	24.48%	5.951%
82	25.737	3850	3855	3862	rVB3	20201	49768	1.50%	0.365%
83	27.041	4071	4077	4088	rVB5	17765	57546	1.74%	0.423%

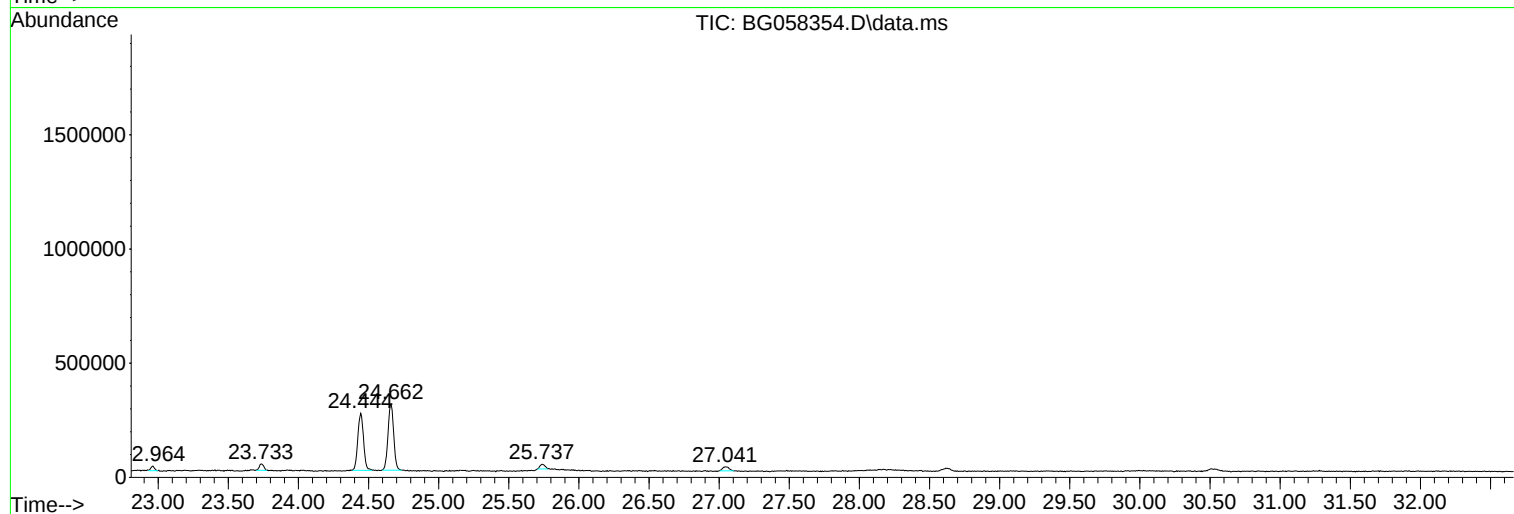
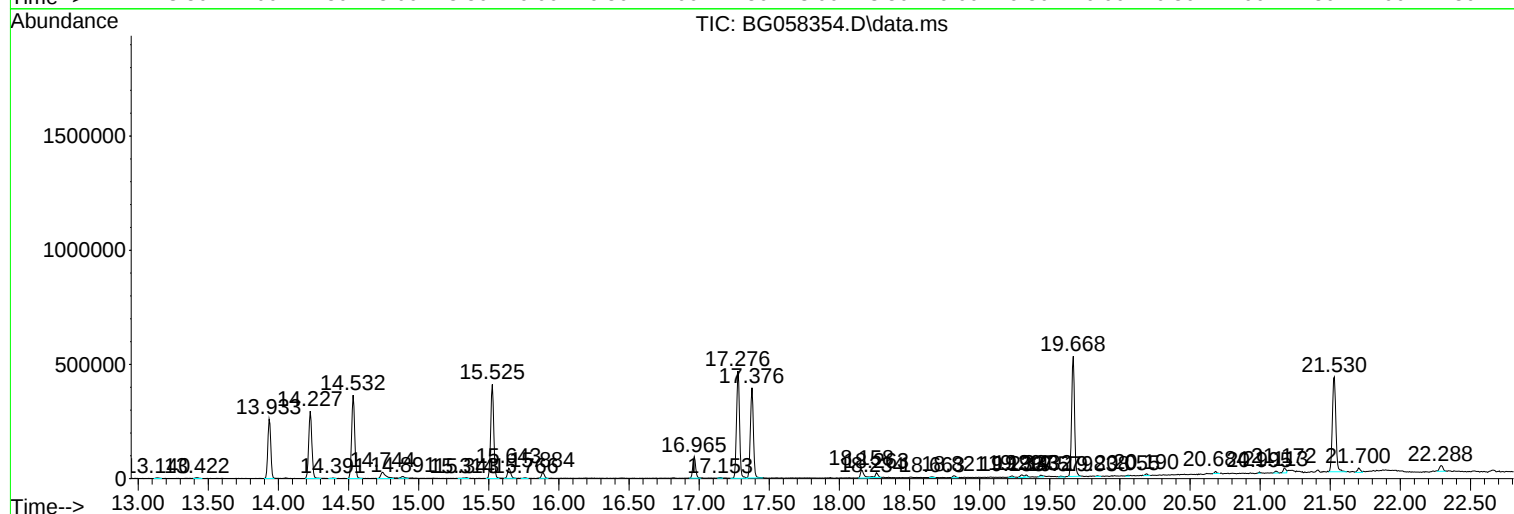
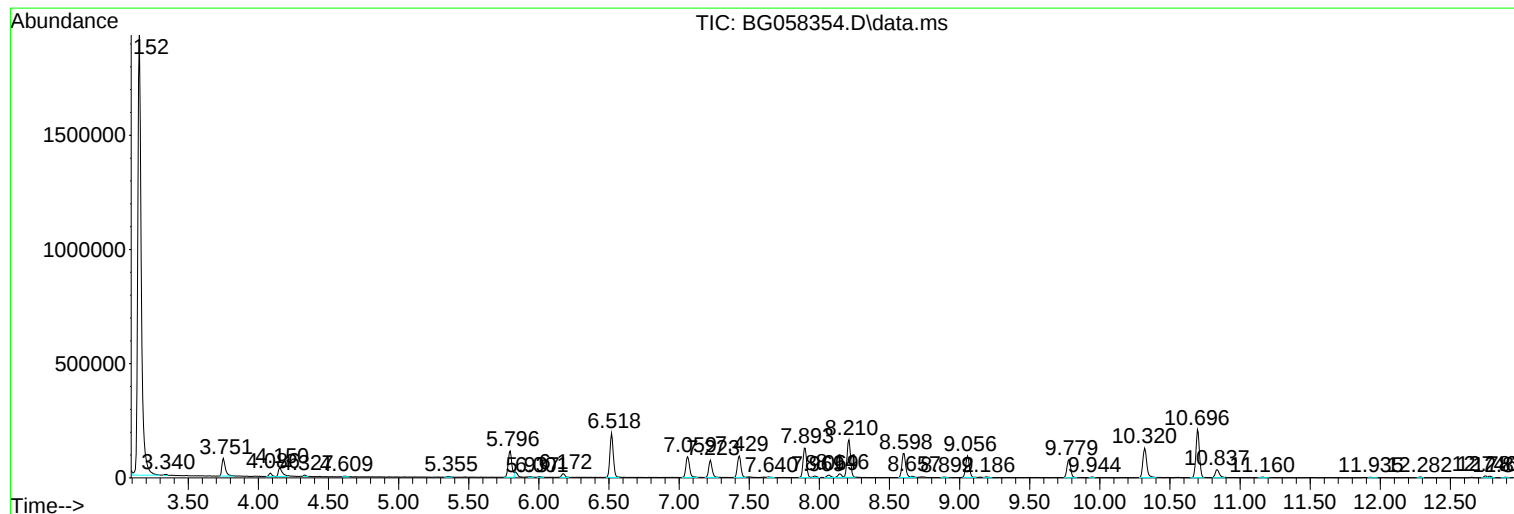
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071723.MA.M
Quant Title : SVOA CALIBRATION

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



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Instrument :
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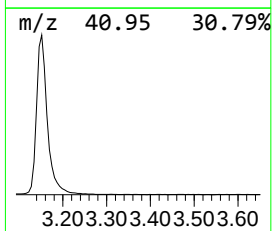
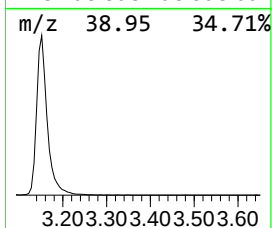
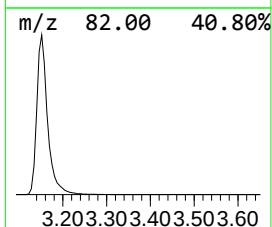
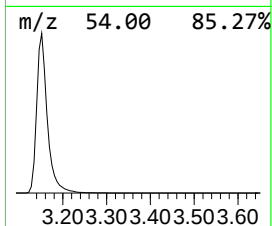
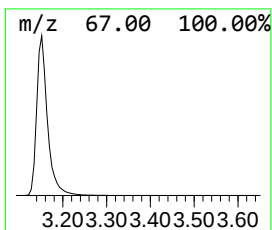
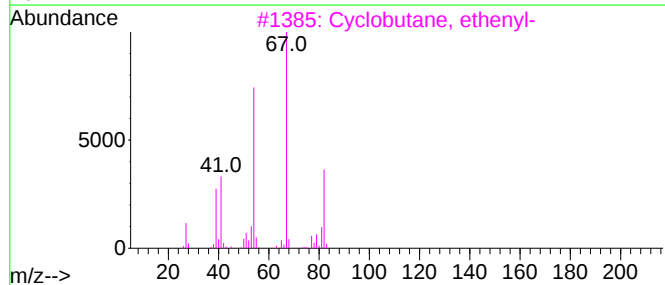
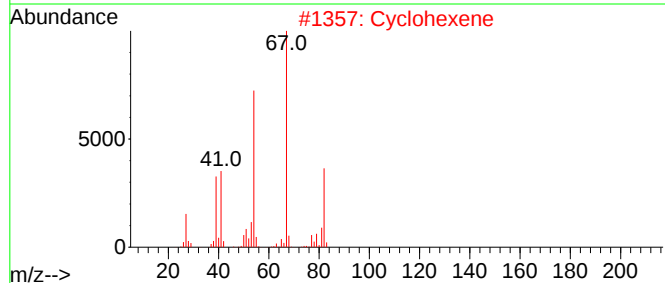
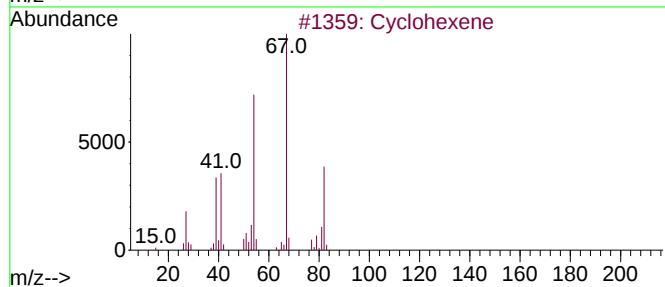
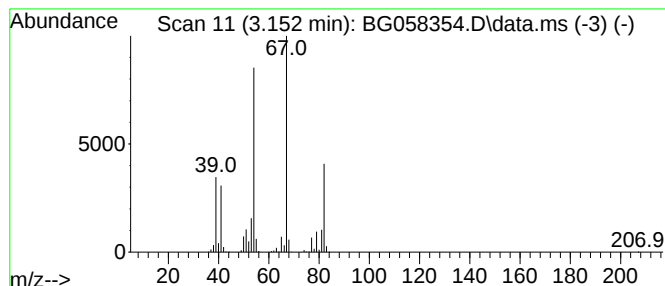
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Cyclohexene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.152	296.80 ng/ul	3310770	1,4-Dichlorobenzene-d4	7.899

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexene	82	C6H10	000110-83-8	95
2			Cyclohexene	82	C6H10	000110-83-8	91
3			Cyclobutane, ethenyl-	82	C6H10	002597-49-1	91
4			Cyclohexene	82	C6H10	000110-83-8	91
5			Cyclohexene	82	C6H10	000110-83-8	90



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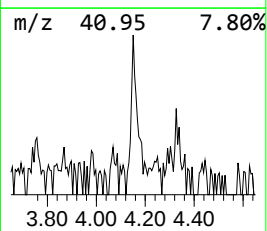
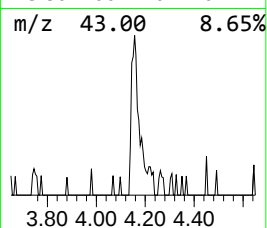
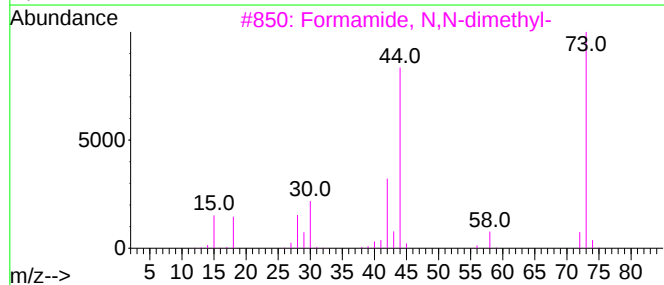
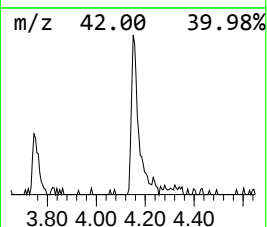
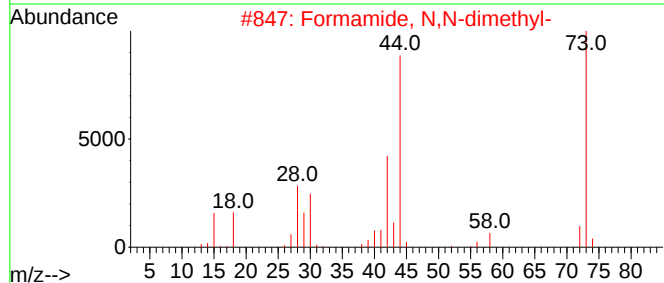
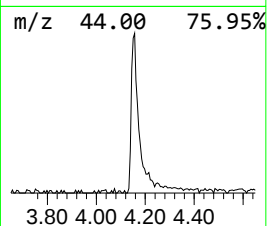
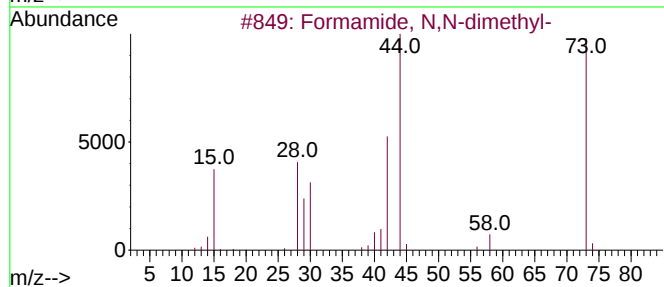
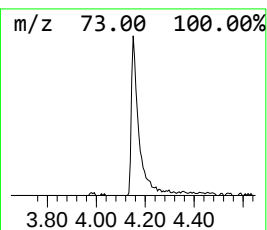
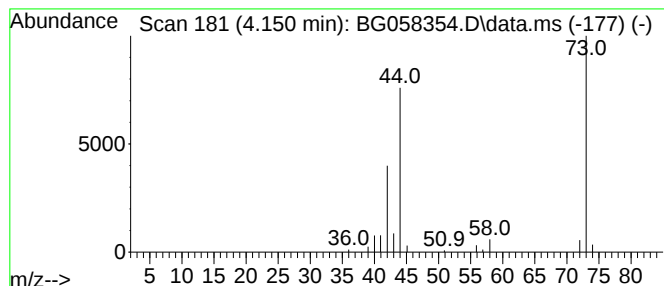
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Formamide, N,N-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.150	6.66 ng/ul	74303	1,4-Dichlorobenzene-d4	7.899

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Formamide, N,N-dimethyl-	73	C3H7NO	000068-12-2	78
2			Formamide, N,N-dimethyl-	73	C3H7NO	000068-12-2	78
3			Formamide, N,N-dimethyl-	73	C3H7NO	000068-12-2	72
4			Formamide, N,N-dimethyl-	73	C3H7NO	000068-12-2	64
5			Glycinamide, N(2)-methyl-	88	C3H8N2O	1000452-55-9	64



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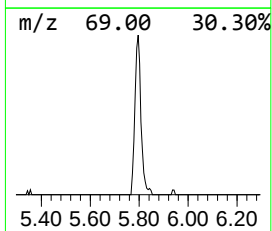
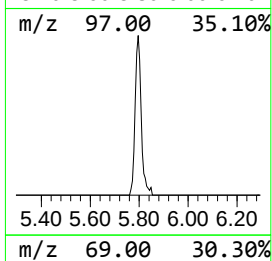
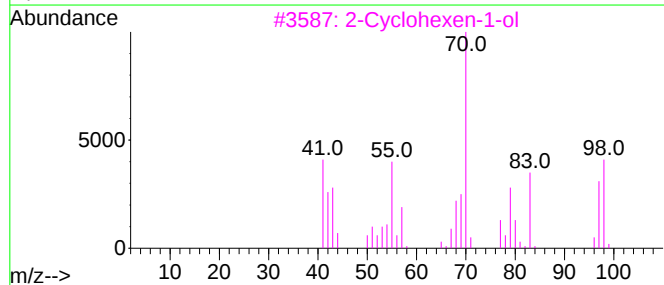
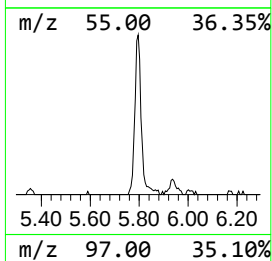
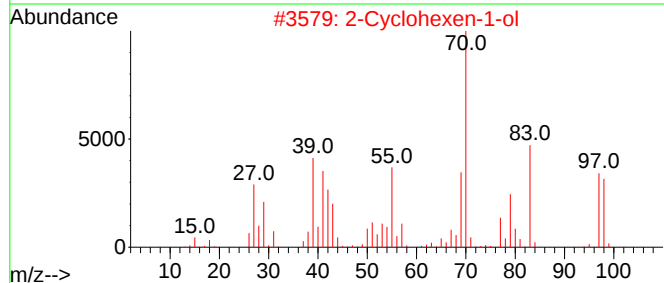
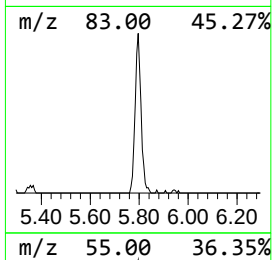
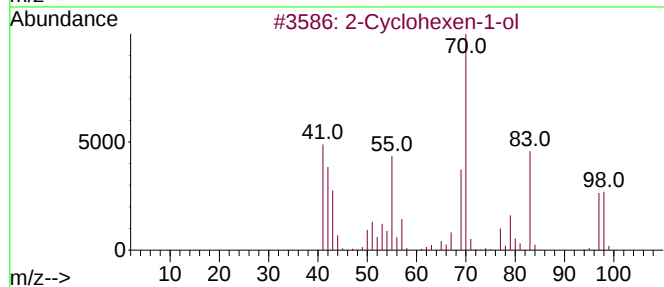
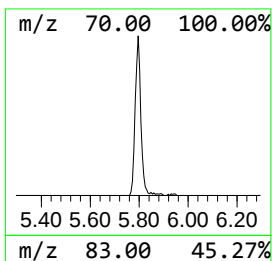
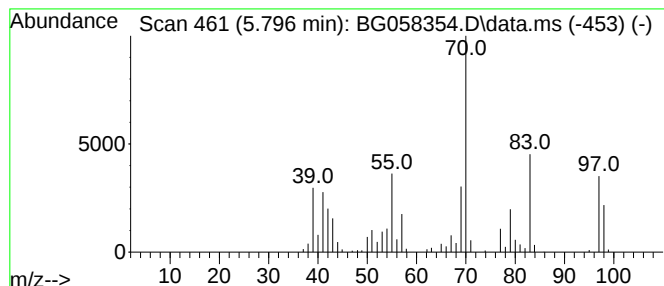
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 4 2-Cyclohexen-1-ol Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.796	19.02 ng/ul	212134	1,4-Dichlorobenzene-d4	7.899

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Cyclohexen-1-ol			98	C6H10O	000822-67-3	93
2	2-Cyclohexen-1-ol			98	C6H10O	000822-67-3	76
3	2-Cyclohexen-1-ol			98	C6H10O	000822-67-3	59
4	2-Methylene cyclopentanol			98	C6H10O	020461-31-8	58
5	Cyclopentane, 1,3-dimethyl-, cis-			98	C7H14	002532-58-3	43



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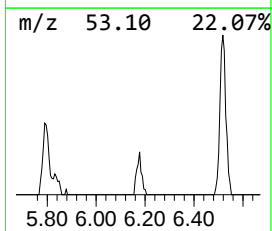
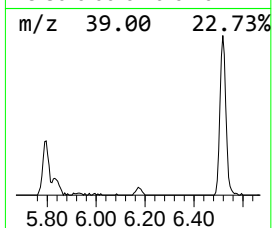
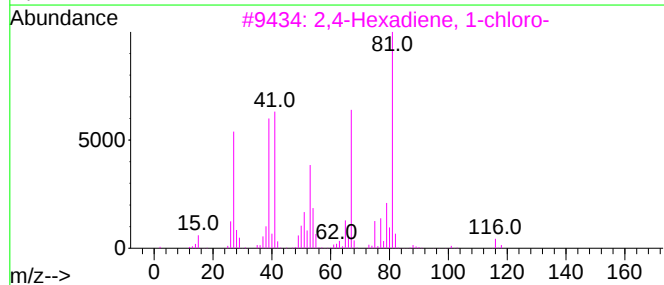
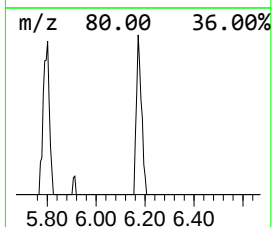
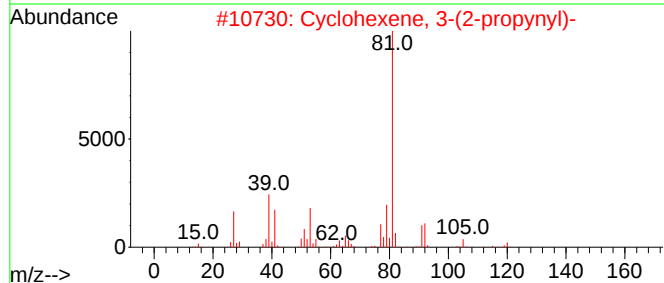
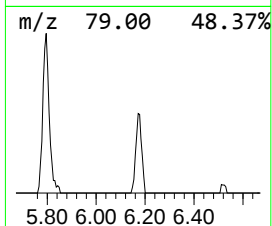
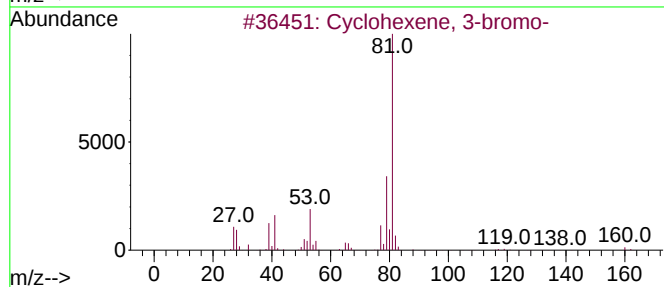
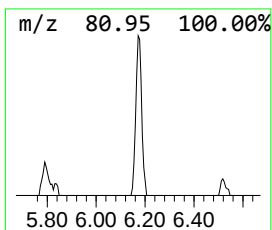
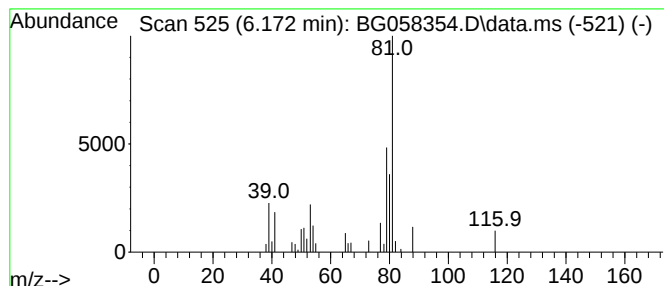
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Cyclohexene, 3-bromo- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.172	2.28 ng/ul	25406	1,4-Dichlorobenzene-d4	7.899

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexene, 3-bromo-	160	C6H9Br	001521-51-3	50
2			Cyclohexene, 3-(2-propynyl)-	120	C9H12	055956-43-9	47
3			2,4-Hexadiene, 1-chloro-	116	C6H9Cl	034632-89-8	43
4			2,3-Dimethyl-1,4-pentadiene	96	C7H12	000758-86-1	38
5			Cyclohexane, 1,2-dichloro-, cis-	152	C6H10Cl2	010498-35-8	36



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071823\
Data File : BG058354.D
Acq On : 20 Jul 2023 9:59
Operator : CG/JU
Sample : 03452-18
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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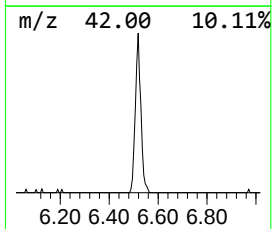
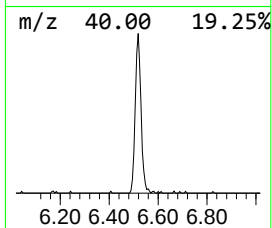
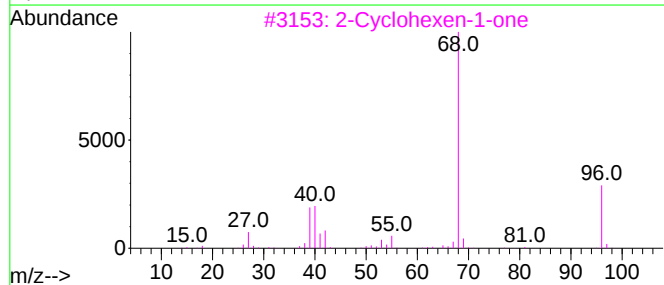
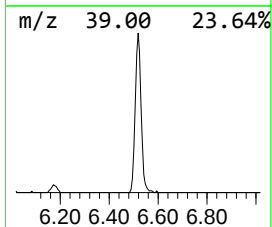
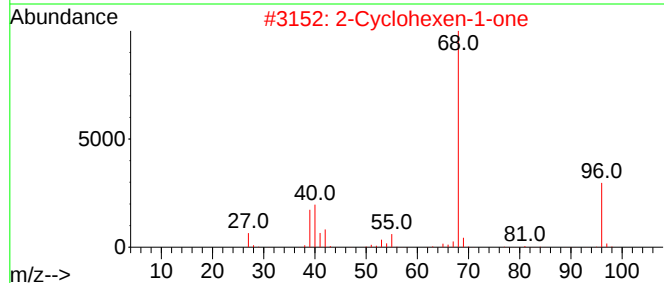
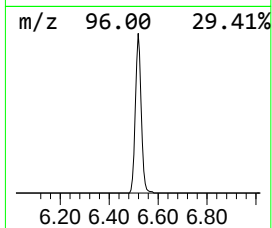
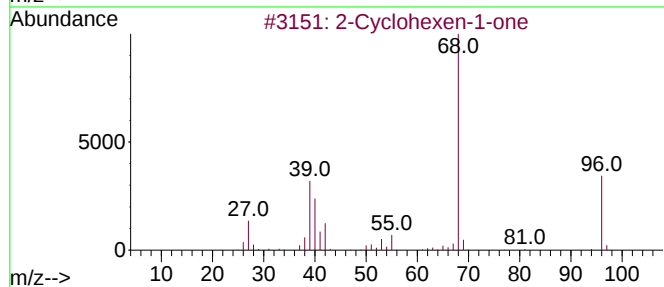
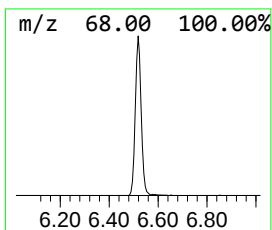
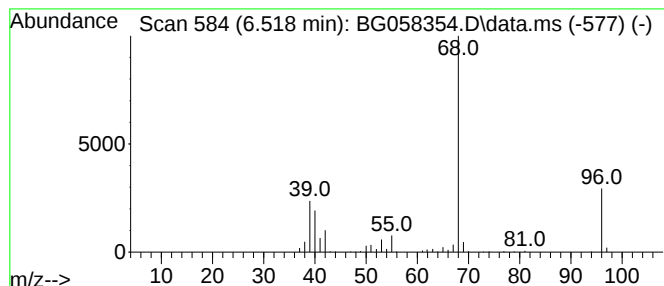
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 2-Cyclohexen-1-one Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.518	28.91 ng/ul	322500	1,4-Dichlorobenzene-d4	7.899

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Cyclohexen-1-one			96	C6H8O	000930-68-7	91
2	2-Cyclohexen-1-one			96	C6H8O	000930-68-7	91
3	2-Cyclohexen-1-one			96	C6H8O	000930-68-7	91
4	2-Cyclohexen-1-one			96	C6H8O	000930-68-7	91
5	7-Oxabicyclo[2.2.1]hept-5-en-2-one			110	C6H6O2	095530-78-2	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071823\
Data File : BG058354.D
Acq On : 20 Jul 2023 9:59
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Sample : 03452-18
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
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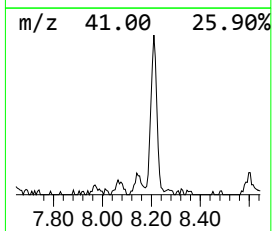
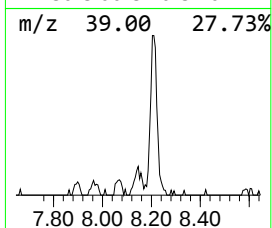
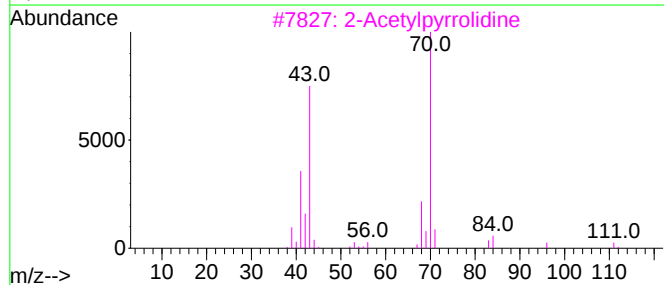
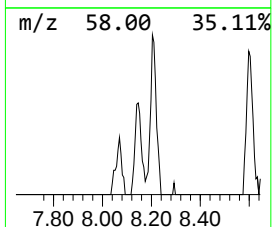
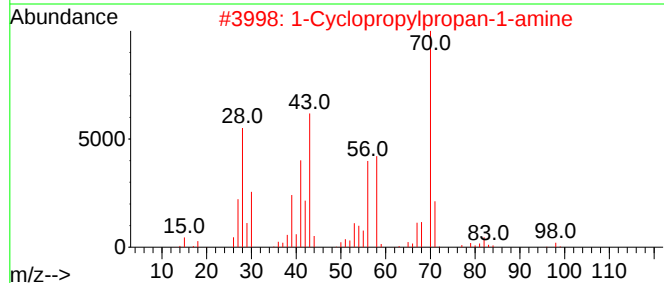
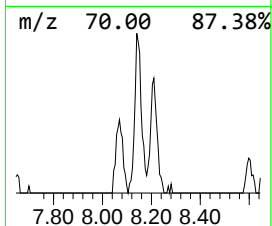
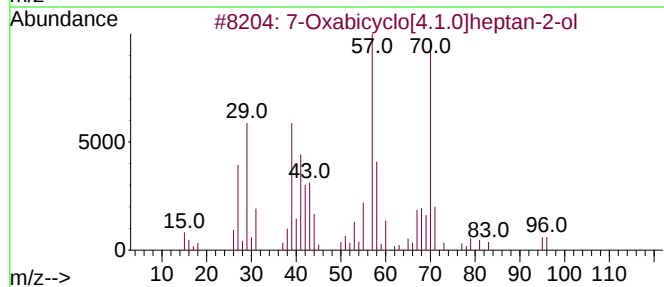
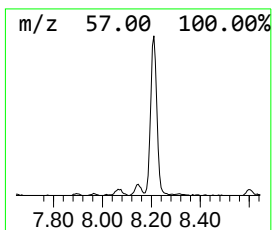
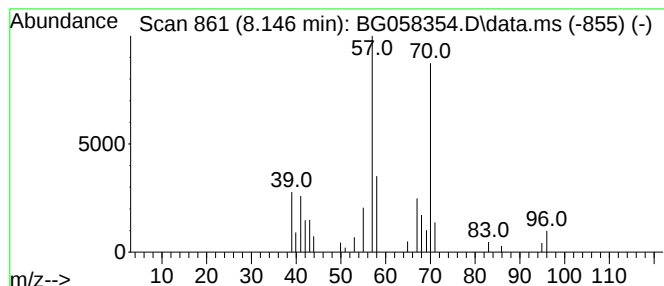
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 7-Oxabicyclo[4.1.0]heptan-2-ol Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.146	2.73 ng/ul	30445	1,4-Dichlorobenzene-d4	7.899

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		7-Oxabicyclo[4.1.0]heptan-2-ol	114	C6H10O2	001192-78-5	50
2		1-Cyclopropylpropan-1-amine	99	C6H13N	1000436-93-8	40
3		2-Acetylpyrrolidine	113	C6H11NO	060026-20-2	28
4		Methacrolein	70	C4H6O	000078-85-3	25
5		2-Heptenol	114	C7H14O	1000429-19-3	10



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071823\
Data File : BG058354.D
Acq On : 20 Jul 2023 9:59
Operator : CG/JU
Sample : 03452-18
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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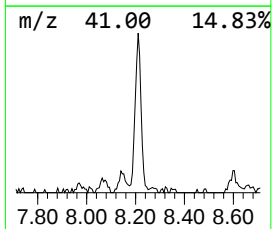
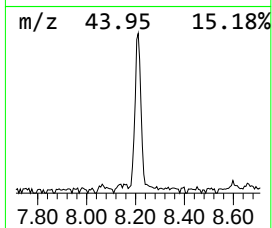
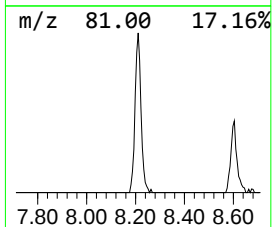
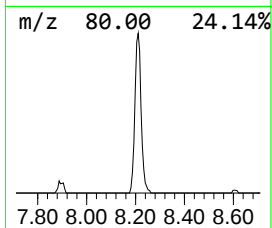
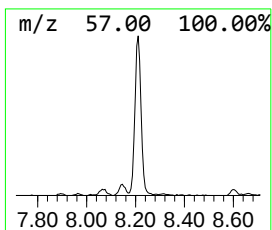
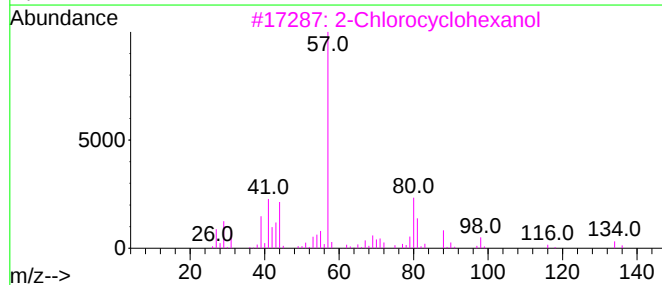
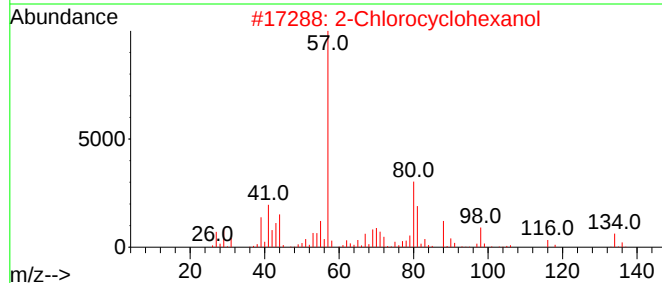
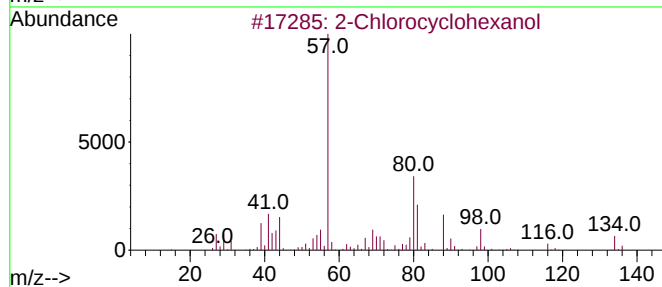
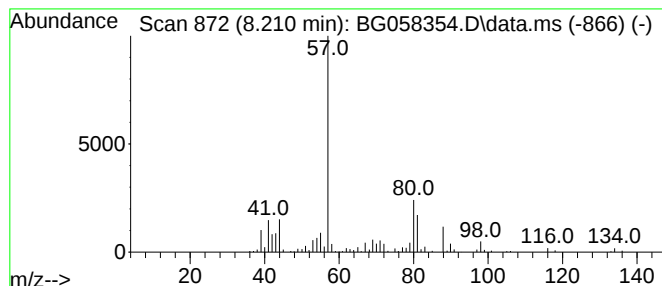
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 2-Chlorocyclohexanol Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.210	25.06 ng/ul	279535	1,4-Dichlorobenzene-d4	7.899

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Chlorocyclohexanol	134	C6H11ClO	001561-86-0	95
2		2-Chlorocyclohexanol	134	C6H11ClO	001561-86-0	94
3		2-Chlorocyclohexanol	134	C6H11ClO	001561-86-0	91
4		2-Chlorocyclohexanol	134	C6H11ClO	001561-86-0	74
5		2-Chlorocyclohexanol	134	C6H11ClO	001561-86-0	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071823\
Data File : BG058354.D
Acq On : 20 Jul 2023 9:59
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Sample : 03452-18
Misc :
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Instrument :
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ClientSampleId :
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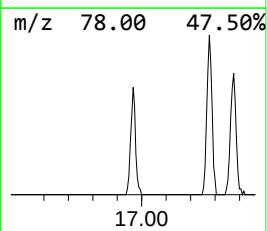
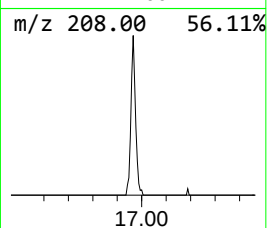
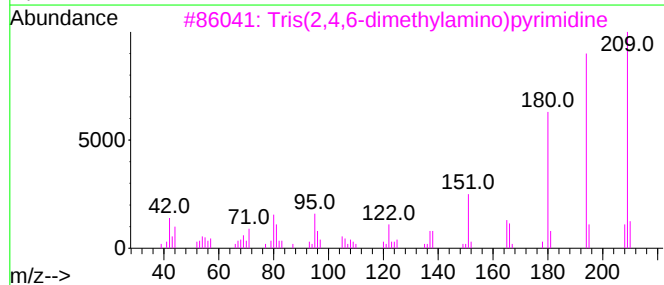
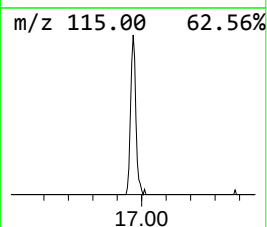
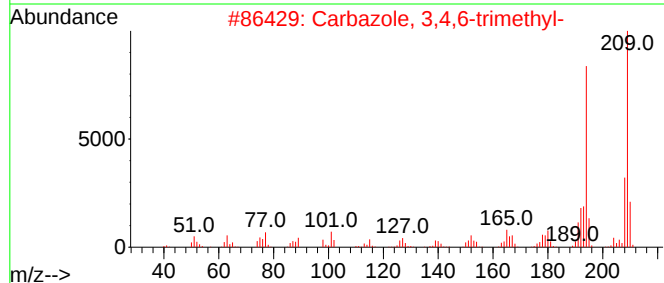
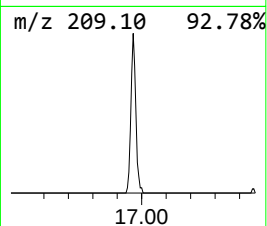
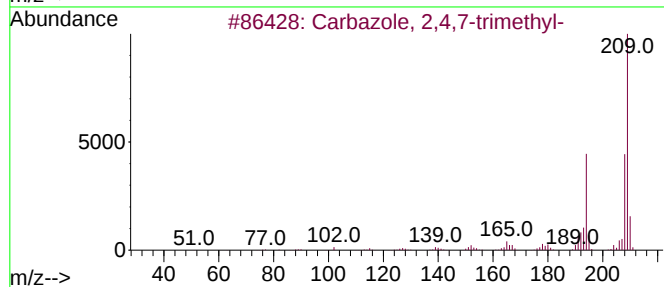
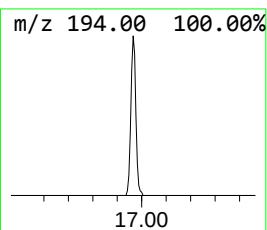
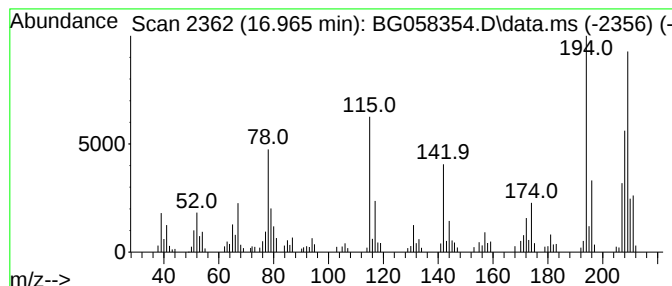
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 unknown-01 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.965	3.69 ng/ul	130636	Phenanthrene-d10	17.276

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Carbazole, 2,4,7-trimethyl-	209	C15H15N	078787-89-0	49
2			Carbazole, 3,4,6-trimethyl-	209	C15H15N	078787-90-3	43
3			Tris(2,4,6-dimethylamino)pyrimidine	209	C10H19N5	043002-24-0	38
4			4-(Dimethylamino)phenol, TMS	209	C11H19NOSi	1000504-21-5	38
5			7-Amino-5,6-dimethoxy-3H-2-benzo...	209	C10H11NO4	500351-80-4	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071823\
Data File : BG058354.D
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Operator : CG/JU
Sample : 03452-18
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071723.MA.M
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
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Formamide, N,N-...	4.150	6.7	ng/ul	74303	1	7.899	223099	20.0
2-Cyclohexen-1-ol	5.796	19.0	ng/ul	212134	1	7.899	223099	20.0
Cyclohexene, 3-...	6.172	2.3	ng/ul	25406	1	7.899	223099	20.0
2-Cyclohexen-1-one	6.518	28.9	ng/ul	322500	1	7.899	223099	20.0
7-Oxabicyclo[4....	8.146	2.7	ng/ul	30445	1	7.899	223099	20.0
2-Chlorocyclohe...	8.210	25.1	ng/ul	279535	1	7.899	223099	20.0
unknown-01	16.965	3.7	ng/ul	130636	4	17.276	707556	20.0