

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122721\
 Data File : BG051800.D
 Acq On : 27 Dec 2021 13:18
 Operator : CG/JU
 Sample : M5201-05
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 P035-TW001-A-01

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-BG121421.M
 Title : SVOA CALIBRATION

Signal : TIC: BG051800.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.602	13	19	25	rVB3	8157	13524	1.19%	0.134%
2	4.030	88	92	99	rBV2	39230	69633	6.15%	0.690%
3	4.083	99	101	105	rVV3	2682	3190	0.28%	0.032%
4	4.383	148	152	156	rBV3	2135	3355	0.30%	0.033%
5	4.465	163	166	170	rVB3	1818	2824	0.25%	0.028%
6	4.894	232	239	247	rBV	121229	195101	17.23%	1.933%
7	5.211	287	293	300	rVB2	10759	17088	1.51%	0.169%
8	5.317	306	311	320	rBV3	1742	4424	0.39%	0.044%
9	5.517	340	345	350	rVB2	5249	7737	0.68%	0.077%
10	6.104	437	445	452	rBV	86443	142999	12.63%	1.417%
11	6.474	502	508	514	rBV3	3300	6259	0.55%	0.062%
12	6.656	534	539	543	rBV2	6351	9715	0.86%	0.096%
13	7.402	659	666	674	rVB2	31197	52898	4.67%	0.524%
14	7.579	688	696	705	rBV	116311	199918	17.65%	1.980%
15	7.796	726	733	742	rBV	127640	223600	19.74%	2.215%
16	8.272	807	814	820	rBV	119736	198729	17.55%	1.969%
17	8.965	926	932	943	rBV	77115	134405	11.87%	1.331%
18	9.147	956	963	970	rBV4	11901	20865	1.84%	0.207%
19	9.441	1003	1013	1022	rBV	165608	292493	25.83%	2.898%
20	10.058	1113	1118	1125	rVV	3119	5054	0.45%	0.050%
21	10.169	1130	1137	1145	rBV	124219	229179	20.24%	2.270%
22	10.463	1180	1187	1190	rBV4	2410	4461	0.39%	0.044%
23	10.715	1222	1230	1239	rBV	184034	330152	29.15%	3.271%
24	10.886	1252	1259	1265	rBV2	34892	60416	5.33%	0.599%
25	11.109	1290	1297	1306	rVB	160132	291351	25.73%	2.886%
26	11.232	1309	1318	1329	rVV	50474	98588	8.71%	0.977%
27	12.302	1493	1500	1508	rVB2	19507	31607	2.79%	0.313%
28	12.678	1556	1564	1571	rVB5	3916	7609	0.67%	0.075%
29	13.059	1621	1629	1639	rBV	49010	81231	7.17%	0.805%
30	13.500	1699	1704	1708	rVV4	3681	5364	0.47%	0.053%
31	14.181	1817	1820	1826	rVB3	6585	9092	0.80%	0.090%
32	14.293	1833	1839	1847	rBV	373286	545982	48.21%	5.409%
33	14.598	1885	1891	1898	rVB	121385	188668	16.66%	1.869%
34	14.904	1936	1943	1951	rBV	266650	418222	36.93%	4.143%
35	15.074	1966	1972	1983	rBV3	31655	61816	5.46%	0.612%
36	15.474	2035	2040	2044	rVB2	1882	3056	0.27%	0.030%

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37	15.897	2105	2112	2122	rBV	568190	858761	75.83%	8.507%
38	15.997	2122	2129	2138	rVV2	145504	217813	19.23%	2.158%
39	16.461	2203	2208	2214	rBV3	5181	8363	0.74%	0.083%
40	16.578	2221	2228	2238	rBV2	87151	151081	13.34%	1.497%
41	17.301	2347	2351	2358	rVB2	16683	25395	2.24%	0.252%
42	17.518	2384	2388	2392	rVB4	3841	5145	0.45%	0.051%
43	17.653	2401	2411	2418	rBV	336580	501537	44.29%	4.968%
44	17.753	2421	2428	2435	rVV2	467568	716782	63.29%	7.101%
45	17.835	2438	2442	2446	rVV3	4009	5665	0.50%	0.056%
46	17.882	2446	2450	2455	rVB5	5873	8461	0.75%	0.084%
47	18.053	2474	2479	2484	rBV3	13785	19998	1.77%	0.198%
48	18.464	2545	2549	2554	rVB2	6522	8675	0.77%	0.086%
49	18.587	2564	2570	2582	rBV	107047	176686	15.60%	1.750%
50	18.869	2613	2618	2624	rBV3	9329	17373	1.53%	0.172%
51	19.022	2640	2644	2648	rVV	10004	13867	1.22%	0.137%
52	19.157	2663	2667	2671	rVV2	11447	14984	1.32%	0.148%
53	19.298	2685	2691	2700	rVB	52990	76635	6.77%	0.759%
54	19.598	2738	2742	2745	rBV3	10060	13461	1.19%	0.133%
55	19.662	2747	2753	2758	rVV	7689	12767	1.13%	0.126%
56	19.885	2787	2791	2796	rBV4	7268	9234	0.82%	0.091%
57	20.026	2809	2815	2822	rBV	786963	1132493	100.00%	11.219%
58	20.120	2829	2831	2834	rBV4	2652	2442	0.22%	0.024%
59	20.338	2865	2868	2871	rBV4	2355	3044	0.27%	0.030%
60	20.802	2944	2947	2949	rBV3	2599	3301	0.29%	0.033%
61	21.049	2985	2989	2990	rBV4	3170	2854	0.25%	0.028%
62	21.101	2995	2998	3000	rVB4	2074	2402	0.21%	0.024%
63	21.272	3020	3027	3030	rBV8	7712	16155	1.43%	0.160%
64	21.571	3076	3078	3080	rBV3	2927	2440	0.22%	0.024%
65	21.601	3082	3083	3086	rBV2	3275	3868	0.34%	0.038%
66	21.871	3127	3129	3131	rVB3	3202	2412	0.21%	0.024%
67	21.971	3139	3146	3152	rBV	300784	513601	45.35%	5.088%
68	22.018	3152	3154	3160	rVB4	13885	21120	1.86%	0.209%
69	22.159	3176	3178	3181	rVV4	2678	3100	0.27%	0.031%
70	22.705	3270	3271	3275	rBV4	2542	3175	0.28%	0.031%
71	22.764	3279	3281	3284	rVV4	2404	3003	0.27%	0.030%
72	22.823	3286	3291	3295	rVB6	6078	11196	0.99%	0.111%
73	23.128	3341	3343	3346	rBV4	3219	4313	0.38%	0.043%
74	23.375	3382	3385	3388	rBV5	2305	3344	0.30%	0.033%
75	23.580	3417	3420	3425	rVB5	7771	11584	1.02%	0.115%
76	23.721	3441	3444	3446	rBV4	3873	5191	0.46%	0.051%

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77	23.997	3489	3491	3493	rBV3	2573	2439	0.22%	0.024%
78	24.056	3498	3501	3502	rBV2	2475	2525	0.22%	0.025%
79	24.467	3566	3571	3577	rBV5	11142	21255	1.88%	0.211%
80	25.219	3692	3699	3715	rVB3	36952	116778	10.31%	1.157%
81	25.472	3729	3742	3758	rVV2	158222	526455	46.49%	5.215%
82	26.265	3875	3877	3881	rBV4	2454	2500	0.22%	0.025%
83	26.347	3889	3891	3894	rBV4	3588	4201	0.37%	0.042%
84	26.776	3957	3964	3974	rVV9	14024	43606	3.85%	0.432%
85	26.841	3974	3975	3978	rVB3	3949	2679	0.24%	0.027%
86	27.029	4003	4007	4008	rBV4	2639	2855	0.25%	0.028%
87	27.240	4041	4043	4046	rBV3	2522	2849	0.25%	0.028%
88	27.587	4100	4102	4105	rBV4	3255	3439	0.30%	0.034%
89	28.104	4188	4190	4192	rBV2	2999	2863	0.25%	0.028%
90	28.292	4206	4222	4241	rBV2	156773	737258	65.10%	7.304%
91	29.337	4397	4400	4401	rBV3	2986	2802	0.25%	0.028%
92	30.095	4527	4529	4530	rVV2	4780	3935	0.35%	0.039%
93	30.119	4531	4533	4539	rVB7	6689	9921	0.88%	0.098%
94	30.324	4565	4568	4569	rBV3	2496	2701	0.24%	0.027%
95	30.359	4572	4574	4577	rBV4	2638	2451	0.22%	0.024%
96	30.841	4654	4656	4658	rBV2	3260	2612	0.23%	0.026%
97	31.552	4776	4777	4781	rBV4	2490	3246	0.29%	0.032%
98	31.640	4790	4792	4796	rBV4	2298	3268	0.29%	0.032%
99	31.934	4841	4842	4846	rBV4	2621	3300	0.29%	0.033%
100	32.310	4902	4906	4907	rVV4	3228	4102	0.36%	0.041%

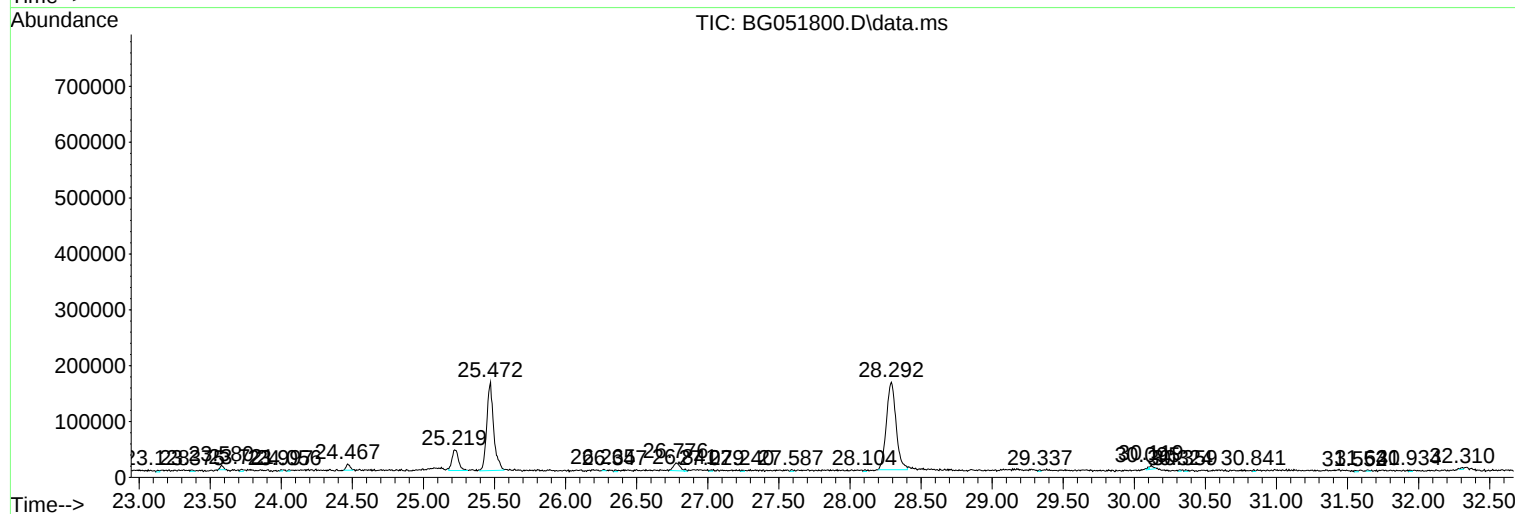
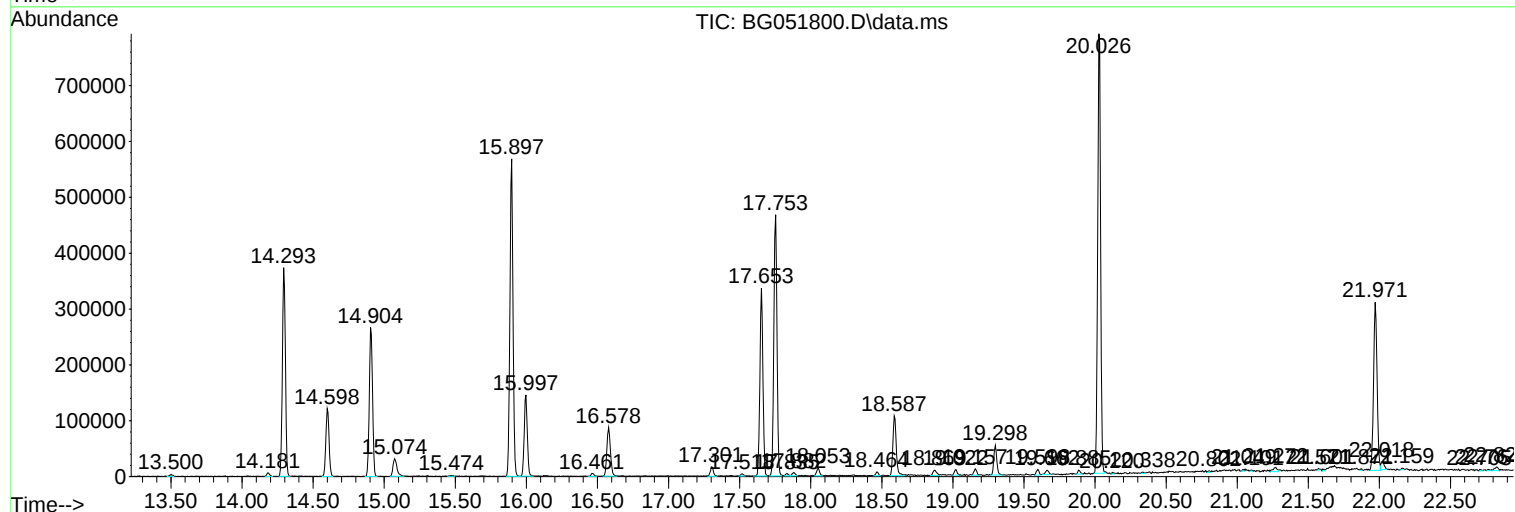
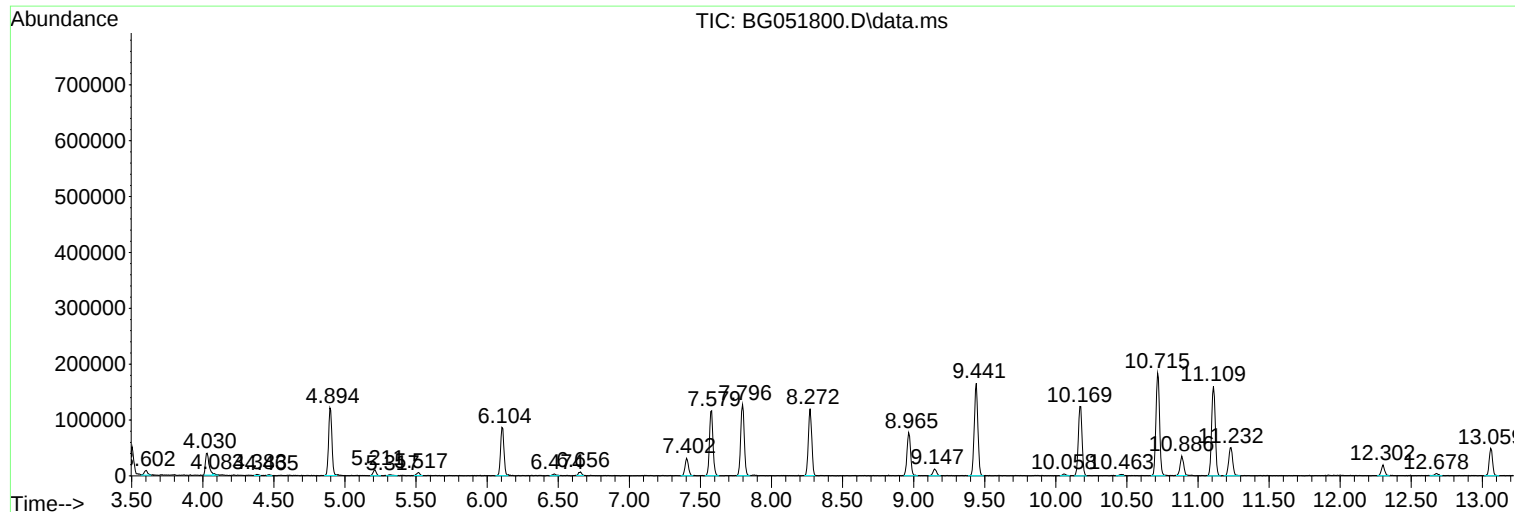
Sum of corrected areas: 10094366

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121421.M
Quant Title : SVOA CALIBRATION

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



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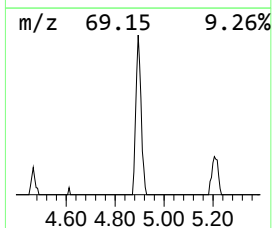
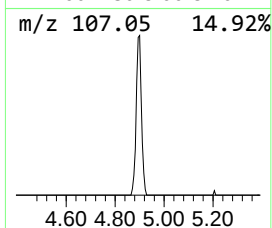
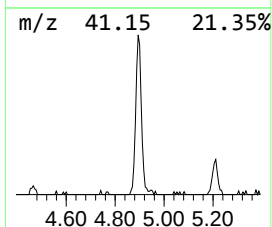
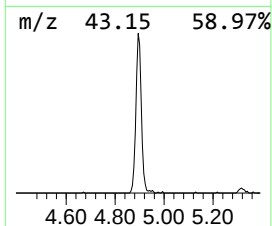
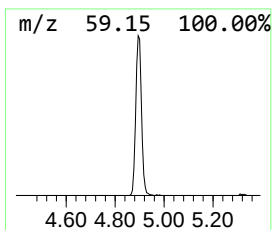
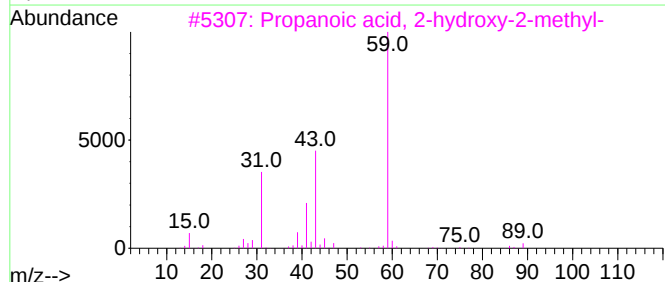
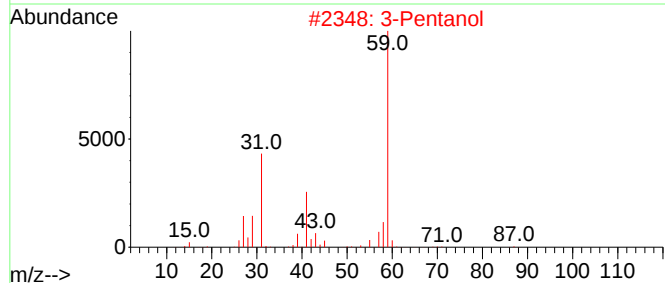
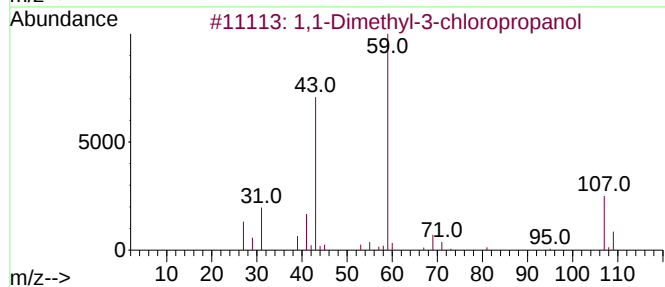
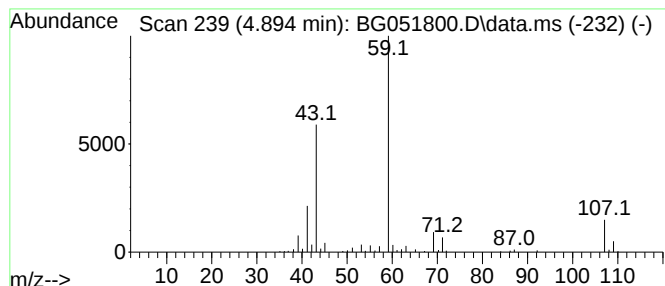
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 1,1-Dimethyl-3-chloropropanol Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.894	19.63 ng/ul	195101	1,4-Dichlorobenzene-d4	8.272

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,1-Dimethyl-3-chloropropanol	122	C5H11ClO	001985-88-2	64
2			3-Pentanol	88	C5H12O	000584-02-1	40
3			Propanoic acid, 2-hydroxy-2-methyl-	104	C4H8O3	000594-61-6	40
4			4-Methyl-2-pentanol, methyl ether	116	C7H16O	1000333-93-5	38
5			4-Methoxy-1-pentene	100	C6H12O	098386-09-5	36



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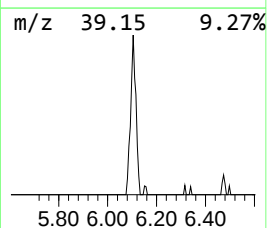
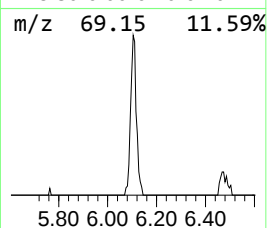
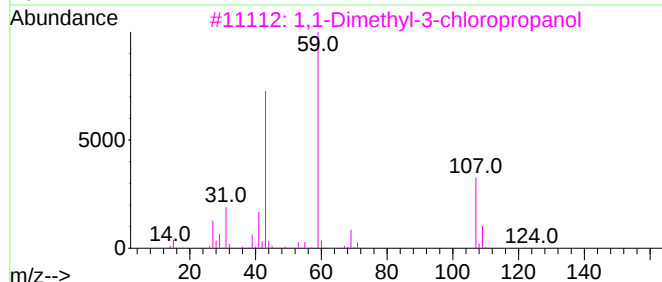
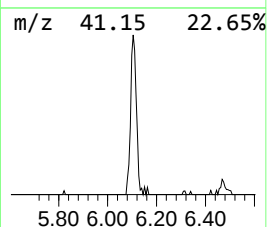
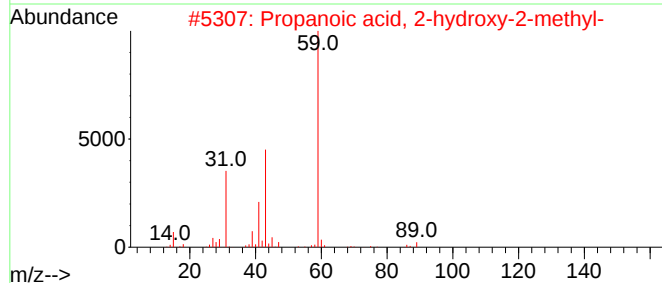
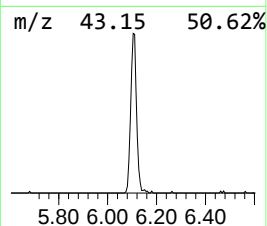
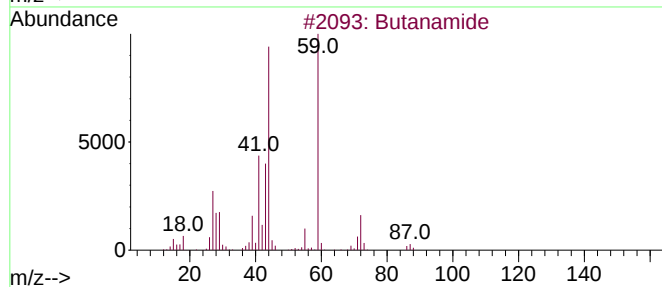
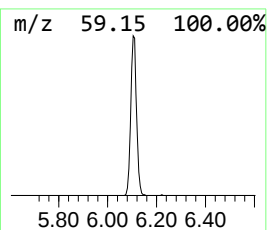
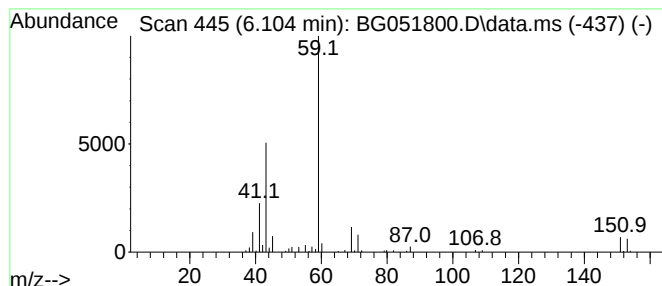
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Butanamide Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.104	14.39 ng/ul	142999	1,4-Dichlorobenzene-d4	8.272

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butanamide	87	C4H9NO	000541-35-5	64
2			Propanoic acid, 2-hydroxy-2-methyl-	104	C4H8O3	000594-61-6	59
3			1,1-Dimethyl-3-chloropropanol	122	C5H11ClO	001985-88-2	56
4			Amylene hydrate	88	C5H12O	000075-85-4	42
5			3-Hydroxy-3-methyl-2-butanone	102	C5H10O2	000115-22-0	42



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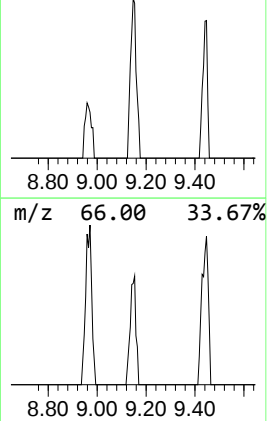
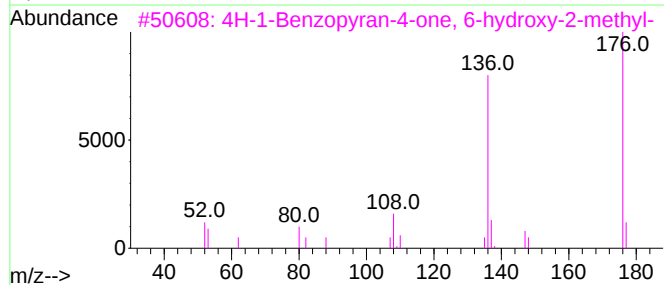
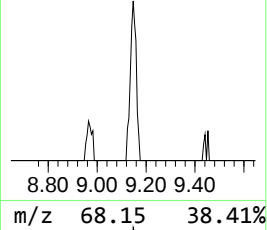
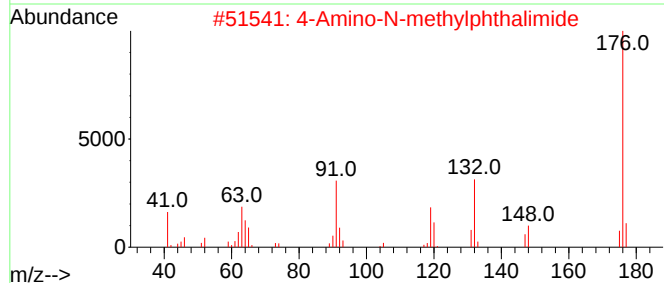
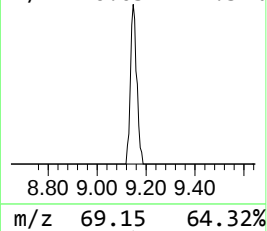
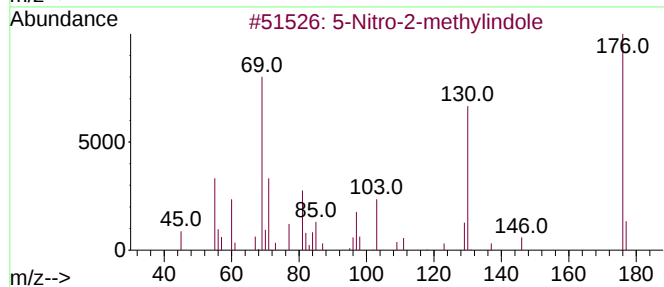
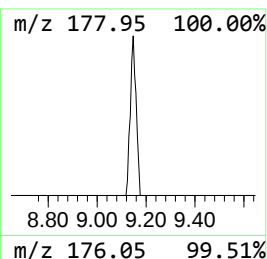
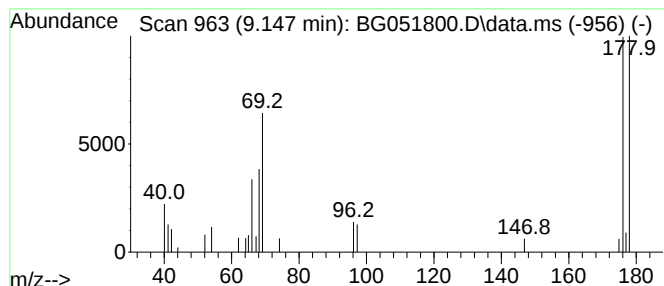
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 unknown-01 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.147	2.10 ng/ul	20865	1,4-Dichlorobenzene-d4	8.272

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			5-Nitro-2-methylindole	176	C9H8N2O2	007570-47-0	14
2			4-Amino-N-methylphthalimide	176	C9H8N2O2	002257-85-4	10
3			4H-1-Benzopyran-4-one, 6-hydroxy...	176	C10H8O3	022105-12-0	10
4			2,4-Pteridinediamine, 6-methyl-	176	C7H8N6	000708-74-7	10
5			m-(Trifluoromethyl)anisole	176	C8H7F3O	000454-90-0	10



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122721\
Data File : BG051800.D
Acq On : 27 Dec 2021 13:18
Operator : CG/JU
Sample : M5201-05
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
P035-TW001-A-01

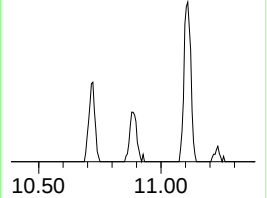
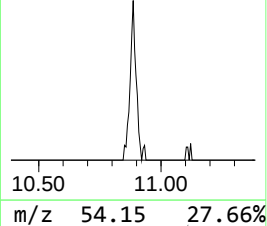
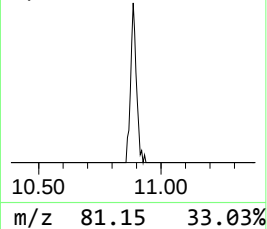
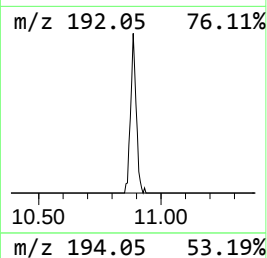
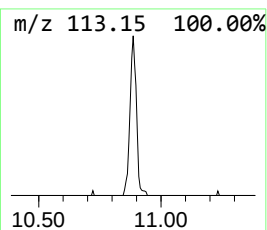
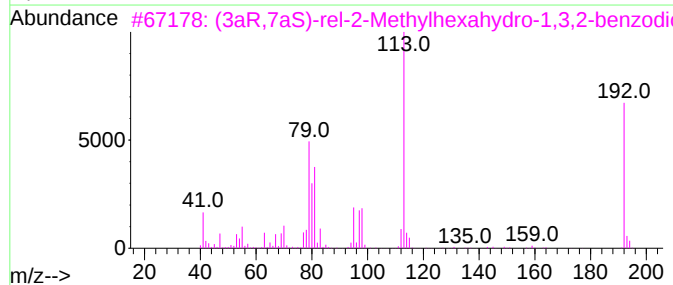
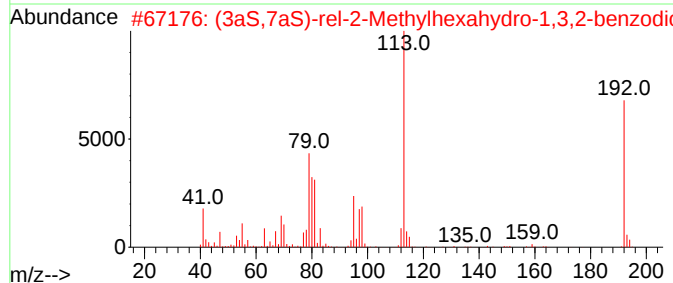
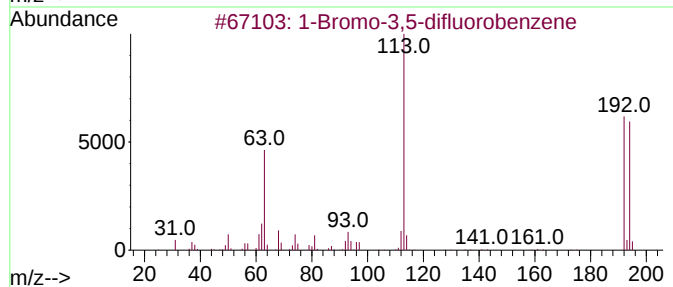
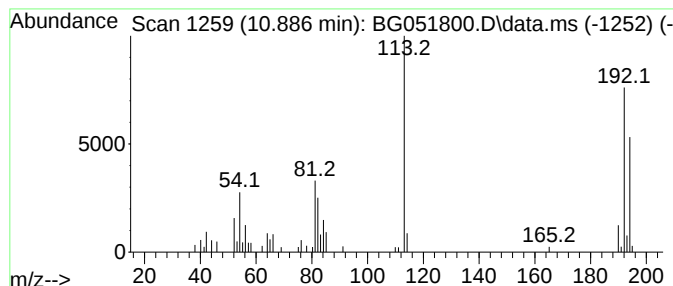
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 1-Bromo-3,5-difluorobenzene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.886	4.15 ng/ul	60416	Naphthalene-d8	11.109

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Bromo-3,5-difluorobenzene	192	C6H3BrF2	000461-96-1	53
2			(3aS,7aS)-rel-2-Methylhexahydro-...	192	C7H13O2PS	1000487-18-9	47
3			(3aR,7aS)-rel-2-Methylhexahydro-...	192	C7H13O2PS	1000487-20-8	47
4			4-Methylamino-2(5H)-furanone	113	C5H7NO2	1000427-45-5	30
5			1-(3-chlorophenyl)-2-(piperazin-...	252	C13H17ClN2O	1000455-11-1	22



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122721\
Data File : BG051800.D
Acq On : 27 Dec 2021 13:18
Operator : CG/JU
Sample : M5201-05
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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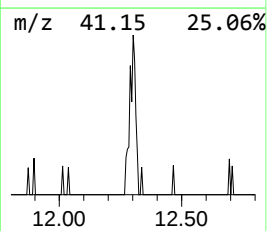
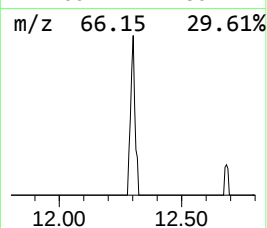
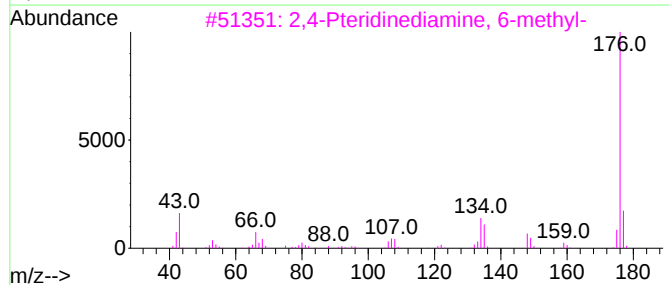
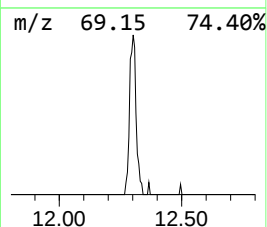
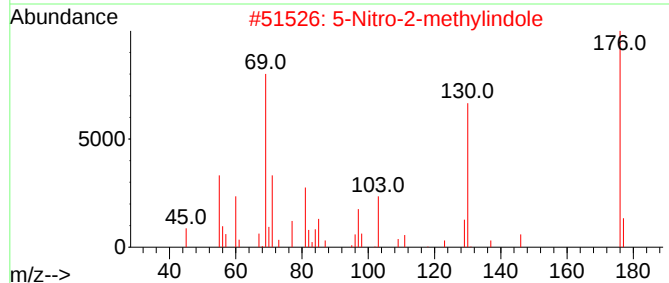
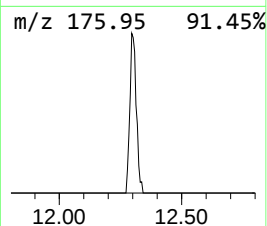
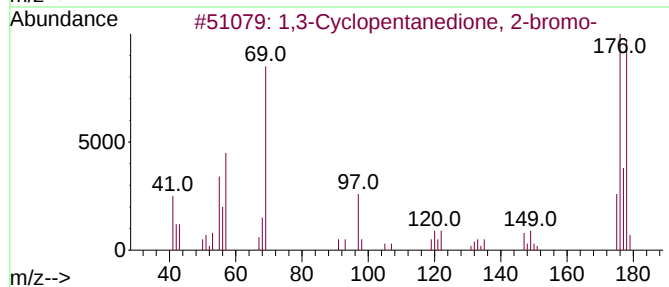
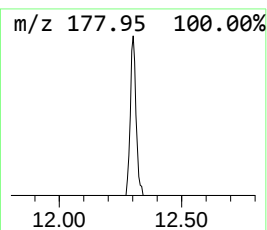
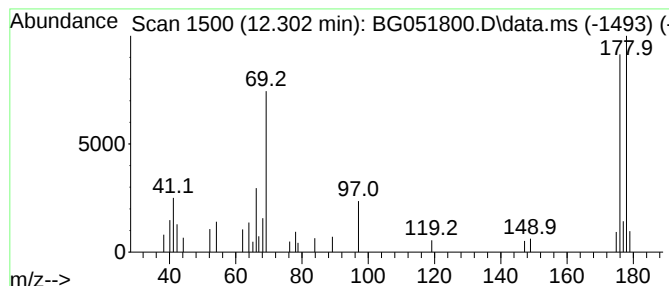
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 unknown-02 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.302	2.17 ng/ul	31607	Naphthalene-d8	11.109

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,3-Cyclopentanedione, 2-bromo-	176	C5H5BrO2	014203-24-8	37
2			5-Nitro-2-methylindole	176	C9H8N2O2	007570-47-0	16
3			2,4-Pteridinediamine, 6-methyl-	176	C7H8N6	000708-74-7	16
4			1,2,4,5-Benzenetetracarbonitrile	178	C10H2N4	000712-74-3	14
5			benzoxazole, 2-methyl-6-nitro-	178	C8H6N2O3	1000404-77-0	14



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122721\
Data File : BG051800.D
Acq On : 27 Dec 2021 13:18
Operator : CG/JU
Sample : M5201-05
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
P035-TW001-A-01

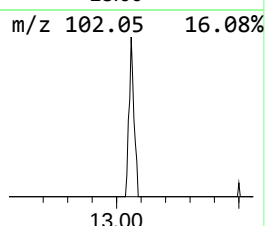
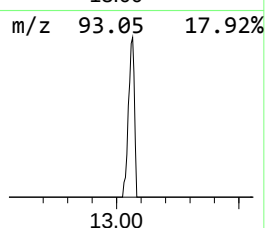
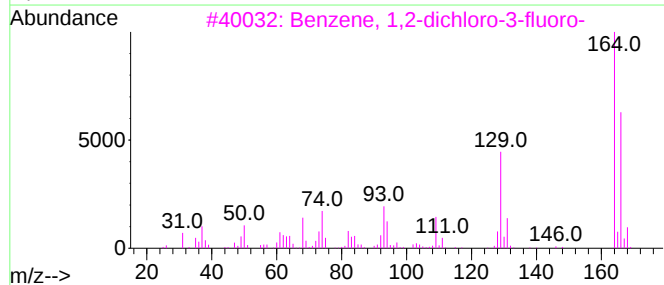
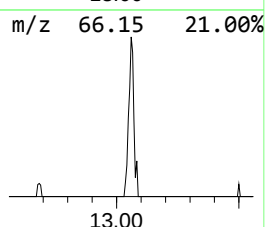
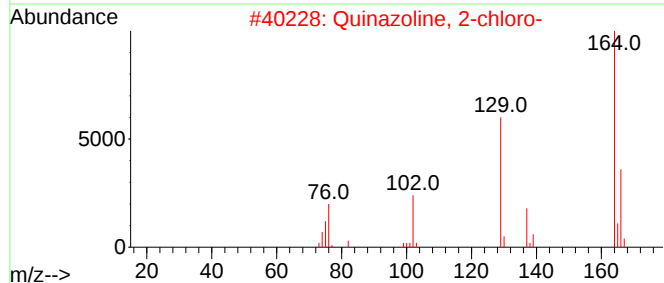
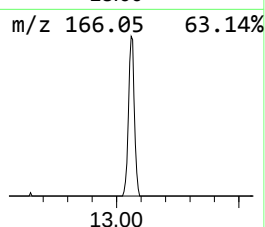
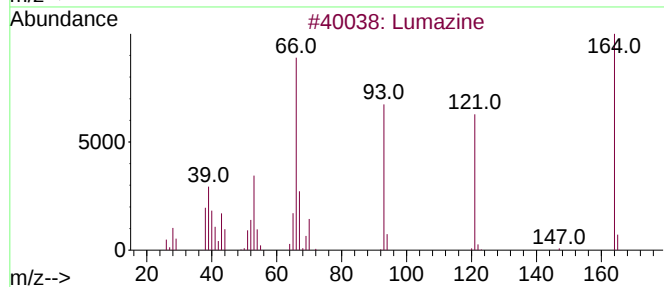
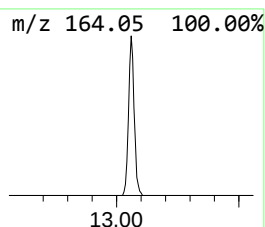
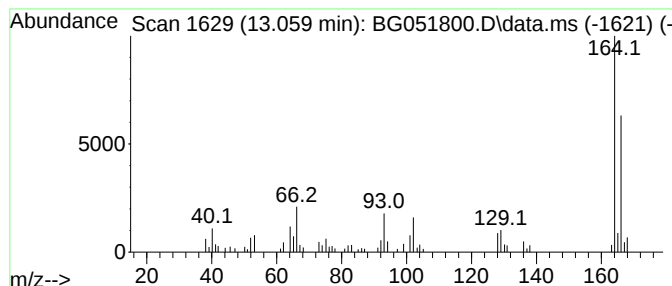
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Lumazine Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.059	3.88 ng/ul	81231	Acenaphthene-d10	14.904

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Lumazine	164	C6H4N4O2	000487-21-8	50
2			Quinazoline, 2-chloro-	164	C8H5ClN2	006141-13-5	45
3			Benzene, 1,2-dichloro-3-fluoro-	164	C6H3Cl2F	036556-50-0	43
4			1,4-Dichloro-2-fluorobenzene	164	C6H3Cl2F	1000342-41-8	43
5			1,3-Dichloro-2-fluorobenzene	164	C6H3Cl2F	002268-05-5	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122721\
Data File : BG051800.D
Acq On : 27 Dec 2021 13:18
Operator : CG/JU
Sample : M5201-05
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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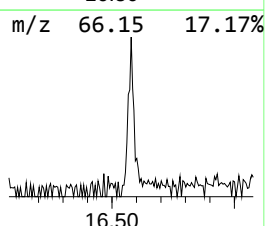
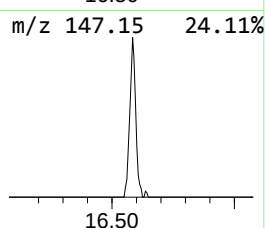
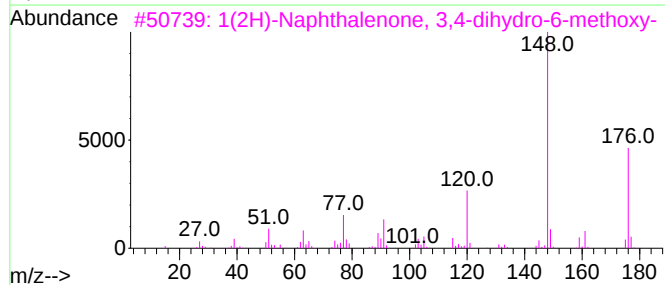
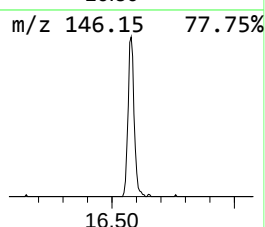
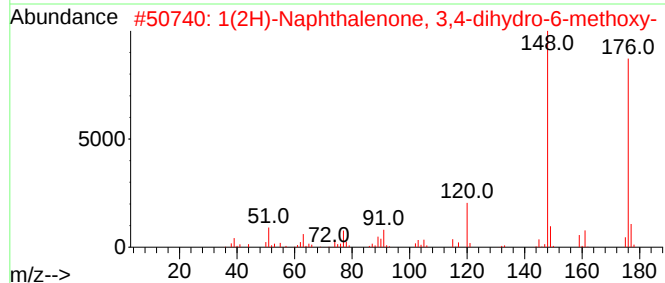
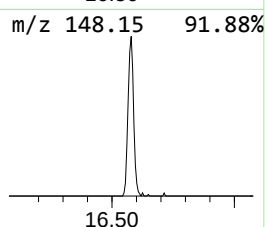
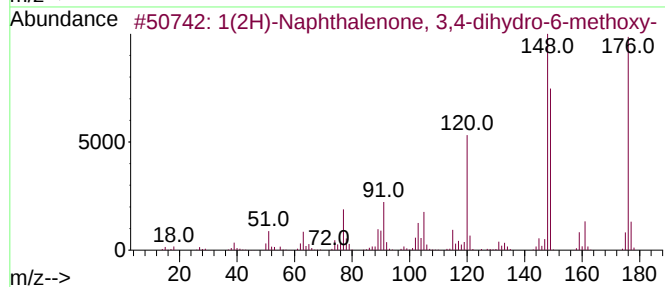
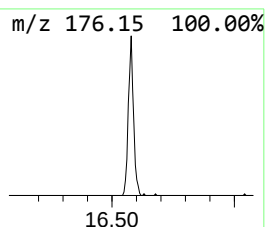
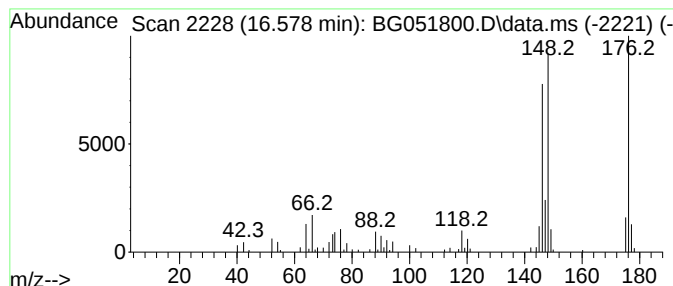
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 unknown-03 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.578	6.02 ng/ul	151081	Phenanthrene-d10	17.653

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1(2H)-Naphthalenone, 3,4-dihydro...	176	C11H12O2	001078-19-9	47		
2	1(2H)-Naphthalenone, 3,4-dihydro...	176	C11H12O2	001078-19-9	47		
3	1(2H)-Naphthalenone, 3,4-dihydro...	176	C11H12O2	001078-19-9	47		
4	1(2H)-Naphthalenone, 3,4-dihydro...	176	C11H12O2	001078-19-9	47		
5	4-Quinazolinol, 2-methyl-, 3-oxide	176	C9H8N2O2	054518-07-9	43		



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122721\
Data File : BG051800.D
Acq On : 27 Dec 2021 13:18
Operator : CG/JU
Sample : M5201-05
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
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ClientSampleId :
P035-TW001-A-01

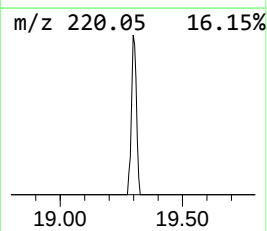
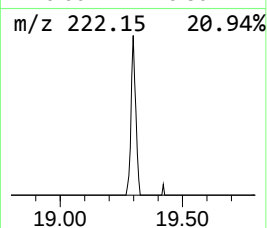
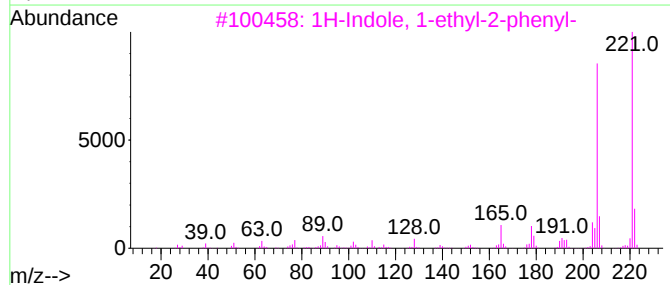
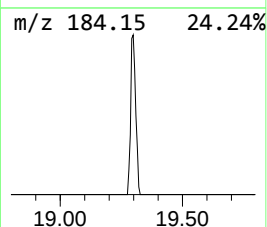
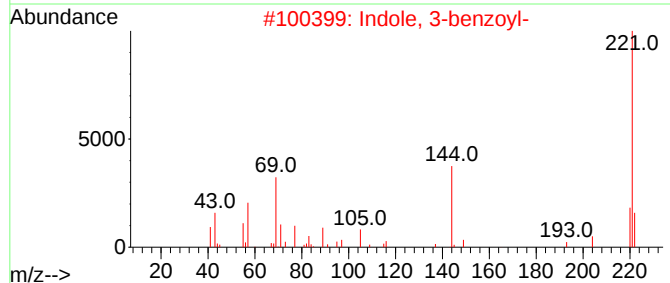
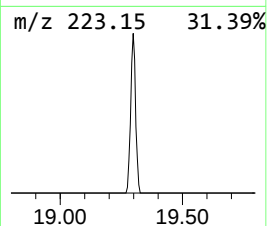
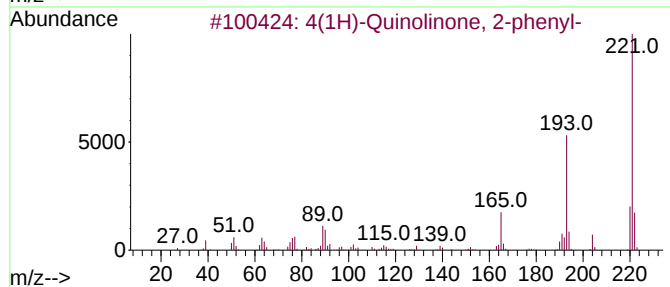
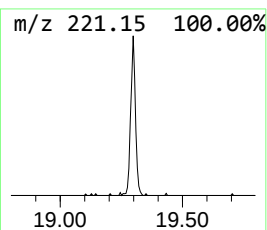
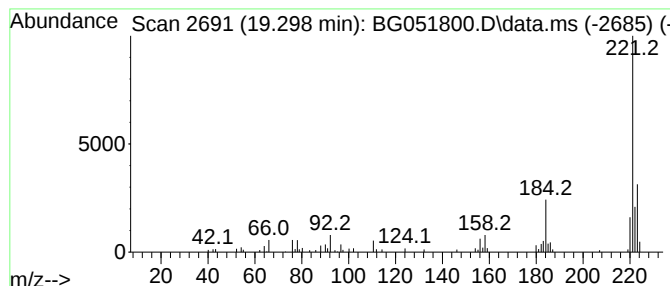
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 unknown-04 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.298	3.06 ng/ul	76635	Phenanthrene-d10	17.653

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	4(1H)-Quinolinone, 2-phenyl-	221	C15H11NO	014802-18-7	43		
2	Indole, 3-benzoyl-	221	C15H11NO	015224-25-6	38		
3	1H-Indole, 1-ethyl-2-phenyl-	221	C16H15N	013228-39-2	38		
4	6-Chloro-2-(trifluoromethyl)-3H-...	221	C7H3ClF3N3	013577-71-4	33		
5	1-(5-Chloro-2-methylphenyl)-2,5-...	221	C11H8ClNO2	058670-26-1	33		



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122721\
Data File : BG051800.D
Acq On : 27 Dec 2021 13:18
Operator : CG/JU
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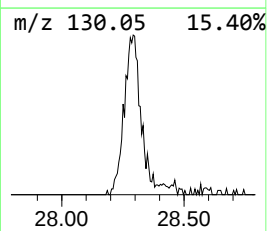
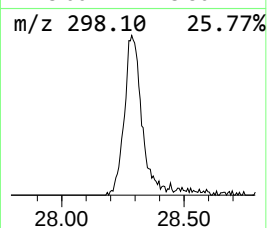
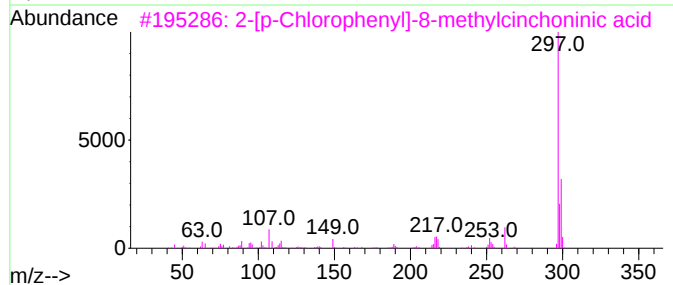
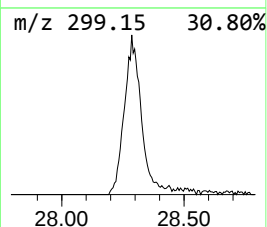
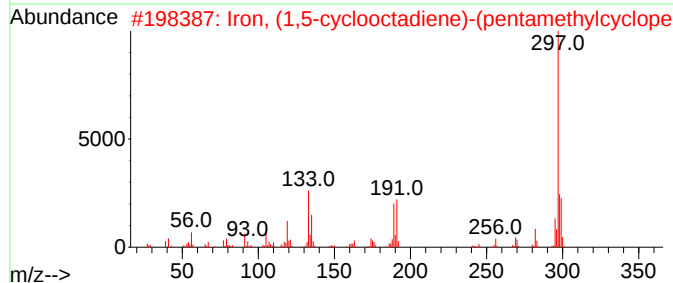
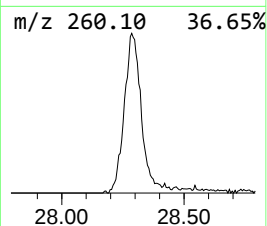
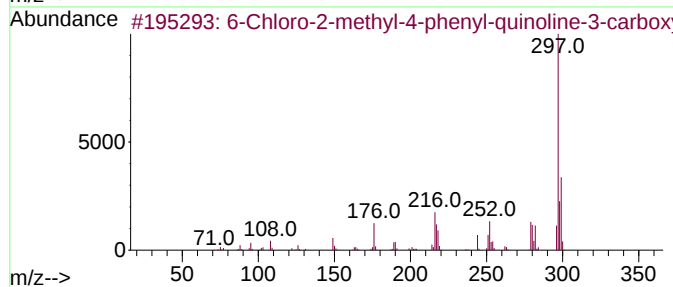
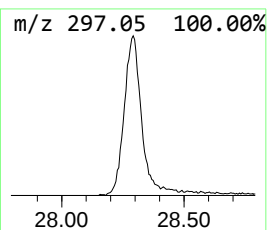
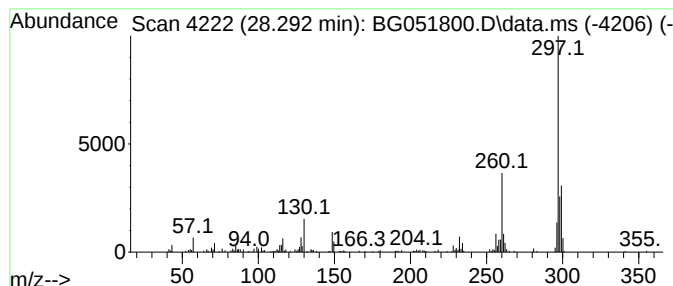
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 6-Chloro-2-methyl-4-phenyl-... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
28.292	28.01 ng/ul	737258	Perylene-d12	25.472

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			6-Chloro-2-methyl-4-phenyl-quinol...	297	C17H12ClNO2	312758-85-3	64
2			Iron, (1,5-cyclooctadiene)-(pent...	299	C18H27Fe	1000161-39-5	64
3			2-[p-Chlorophenyl]-8-methylcinch...	297	C17H12ClNO2	107027-43-0	52
4			6-(Adamantan-1-yl)-2-chloropyrid...	297	C17H16ClN3	1000410-36-2	50
5			trans-4'-Dimethylamino-4-(methyl...	297	C18H19NOS	126443-24-1	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122721\
 Data File : BG051800.D
 Acq On : 27 Dec 2021 13:18
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 ALS Vial : 7 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121421.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
1,1-Dimethyl-3-...	4.894	19.6	ng/ul	195101	1	8.272	198729	20.0
Butanamide	6.104	14.4	ng/ul	142999	1	8.272	198729	20.0
unknown-01	9.147	2.1	ng/ul	20865	1	8.272	198729	20.0
1-Bromo-3,5-dif...	10.886	4.2	ng/ul	60416	2	11.109	291351	20.0
unknown-02	12.302	2.2	ng/ul	31607	2	11.109	291351	20.0
Lumazine	13.059	3.9	ng/ul	81231	3	14.904	418222	20.0
unknown-03	16.578	6.0	ng/ul	151081	4	17.653	501537	20.0
unknown-04	19.298	3.1	ng/ul	76635	4	17.653	501537	20.0
6-Chloro-2-meth...	28.292	28.0	ng/ul	737258	6	25.472	526455	20.0