

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121421\
 Data File : BG051505.D
 Acq On : 14 Dec 2021 15:46
 Operator : CG/JU
 Sample : M5013-05
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 DBLS7

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: BG051505.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.608	14	20	26	rVV	6725	11418	0.56%	0.033%
2	3.966	76	81	82	rBV4	2381	2787	0.14%	0.008%
3	4.043	89	94	104	rBV	92897	163894	8.05%	0.478%
4	4.395	148	154	160	rVV2	3814	7242	0.36%	0.021%
5	4.659	193	199	203	rBV3	4706	7644	0.38%	0.022%
6	4.718	206	209	213	rBV4	4463	6164	0.30%	0.018%
7	4.930	240	245	247	rBV5	2461	3674	0.18%	0.011%
8	5.335	308	314	322	rVB	26192	39166	1.92%	0.114%
9	5.552	345	351	357	rVB2	18601	32218	1.58%	0.094%
10	6.439	496	502	508	rBV3	7869	14794	0.73%	0.043%
11	7.274	640	644	650	rVB4	5216	8938	0.44%	0.026%
12	7.326	650	653	658	rBV3	1282	3021	0.15%	0.009%
13	7.409	660	667	677	rVB	180473	328576	16.14%	0.957%
14	7.491	677	681	682	rBV2	1782	2498	0.12%	0.007%
15	7.514	682	685	688	rVB4	2011	3153	0.15%	0.009%
16	7.591	691	698	707	rVV	105235	175175	8.61%	0.510%
17	7.808	726	735	737	rBV	151239	247508	12.16%	0.721%
18	7.837	737	740	749	rVB	181604	311361	15.30%	0.907%
19	7.908	749	752	757	rBV2	1512	2650	0.13%	0.008%
20	8.284	809	816	823	rBV	108777	186830	9.18%	0.544%
21	8.531	852	858	863	rBV4	2363	4651	0.23%	0.014%
22	8.813	896	906	913	rVV2	181836	451488	22.18%	1.316%
23	8.865	913	915	921	rVB3	17463	25196	1.24%	0.073%
24	8.977	926	934	944	rBV	188771	323137	15.88%	0.942%
25	9.112	949	957	964	rBV	227599	403270	19.81%	1.175%
26	9.388	997	1004	1009	rBV	86419	150244	7.38%	0.438%
27	9.453	1009	1015	1018	rVV	141328	257683	12.66%	0.751%
28	9.494	1018	1022	1033	rVB	257236	450220	22.12%	1.312%
29	9.770	1062	1069	1077	rBV	340689	555258	27.28%	1.618%
30	10.187	1132	1140	1141	rVV	115043	187975	9.24%	0.548%
31	10.217	1141	1145	1152	rVV	185629	345183	16.96%	1.006%
32	10.281	1152	1156	1163	rVB4	11838	24103	1.18%	0.070%
33	10.728	1224	1232	1244	rVV2	204218	590450	29.01%	1.720%
34	10.974	1269	1274	1280	rVV3	13632	23705	1.16%	0.069%
35	11.121	1288	1299	1304	rVV	146661	278826	13.70%	0.812%
36	11.174	1304	1308	1314	rVV	119759	208071	10.22%	0.606%

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Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121421.M
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37	11.239	1314	1319	1331	rVB	76135	139285	6.84%	0.406%
38	11.820	1412	1418	1425	rVV	11423	16639	0.82%	0.048%
39	12.367	1503	1511	1519	rBV	336298	560220	27.52%	1.632%
40	12.625	1548	1555	1563	rBV	251214	416437	20.46%	1.213%
41	12.684	1563	1565	1569	rVB3	2470	3332	0.16%	0.010%
42	12.884	1595	1599	1601	rBV3	3916	4614	0.23%	0.013%
43	12.978	1608	1615	1625	rVB	448730	753152	37.00%	2.195%
44	13.124	1633	1640	1646	rBV2	177765	290818	14.29%	0.847%
45	13.354	1672	1679	1684	rBV	241809	397961	19.55%	1.160%
46	13.753	1741	1747	1754	rVV	487095	760784	37.38%	2.217%
47	13.988	1779	1787	1795	rBV	319286	512234	25.17%	1.493%
48	14.105	1802	1807	1813	rVB	10047	15334	0.75%	0.045%
49	14.311	1835	1842	1848	rBV	392034	558864	27.46%	1.628%
50	14.399	1850	1857	1863	rVV	128568	198026	9.73%	0.577%
51	14.470	1863	1869	1877	rVB	285964	447220	21.97%	1.303%
52	14.617	1887	1894	1896	rVV	404720	717546	35.25%	2.091%
53	14.640	1896	1898	1907	rVB	576947	806086	39.60%	2.349%
54	14.916	1936	1945	1951	rBV	256661	403613	19.83%	1.176%
55	14.981	1951	1956	1962	rVB	61114	93786	4.61%	0.273%
56	15.092	1966	1975	1983	rBV2	475929	923653	45.38%	2.691%
57	15.257	1994	2003	2008	rBV	363713	540256	26.54%	1.574%
58	15.316	2008	2013	2019	rVB	247944	389292	19.13%	1.134%
59	15.545	2045	2052	2062	rVB	84943	139952	6.88%	0.408%
60	15.639	2065	2068	2069	rBV3	1916	2445	0.12%	0.007%
61	15.715	2075	2081	2087	rVB	632342	861148	42.31%	2.509%
62	15.903	2107	2113	2119	rBV	521773	806428	39.62%	2.350%
63	15.962	2119	2123	2126	rVV	125806	187430	9.21%	0.546%
64	16.003	2126	2130	2136	rVB	170229	252368	12.40%	0.735%
65	16.156	2145	2156	2162	rBV	774449	1190412	58.49%	3.469%
66	16.391	2191	2196	2205	rBV2	60903	98749	4.85%	0.288%
67	16.473	2207	2210	2211	rVV3	3146	2623	0.13%	0.008%
68	16.532	2217	2220	2225	rVV	12358	18931	0.93%	0.055%
69	16.614	2232	2234	2237	rBV3	1975	2435	0.12%	0.007%
70	16.655	2240	2241	2244	rVB2	2292	2330	0.11%	0.007%
71	16.690	2244	2247	2249	rBV3	2536	3223	0.16%	0.009%
72	16.767	2257	2260	2262	rBV3	4187	4015	0.20%	0.012%
73	16.837	2267	2272	2275	rVV2	12392	16916	0.83%	0.049%
74	16.902	2275	2283	2289	rVV	159372	231058	11.35%	0.673%
75	16.972	2289	2295	2303	rVB	515942	791088	38.87%	2.305%
76	17.307	2346	2352	2362	rBV	416826	615979	30.26%	1.795%

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77	17.601	2398	2402	2405	rBV3	2225	3329	0.16%	0.010%
78	17.665	2406	2413	2416	rVV	368988	569146	27.96%	1.658%
79	17.706	2416	2420	2425	rVV	593254	886937	43.58%	2.584%
80	17.765	2425	2430	2440	rVB	603182	931235	45.75%	2.713%
81	18.000	2468	2470	2473	rVB4	3218	2870	0.14%	0.008%
82	18.276	2512	2517	2521	rBV3	26021	34070	1.67%	0.099%
83	18.611	2569	2574	2580	rVB	748470	941340	46.25%	2.743%
84	19.598	2740	2742	2748	rVB4	9175	13972	0.69%	0.041%
85	19.669	2752	2754	2758	rVB4	9963	11928	0.59%	0.035%
86	19.898	2788	2793	2798	rBV2	80090	115113	5.66%	0.335%
87	20.039	2811	2817	2820	rBV	787766	1292540	63.50%	3.766%
88	20.256	2850	2854	2863	rBV	131301	178001	8.75%	0.519%
89	21.061	2988	2991	2997	rVB	65046	83852	4.12%	0.244%
90	21.860	3124	3127	3133	rVB2	43726	69869	3.43%	0.204%
91	21.965	3139	3145	3156	rVB2	889848	2035347	100.00%	5.931%
92	22.347	3207	3210	3216	rVB7	45434	75024	3.69%	0.219%
93	22.576	3242	3249	3256	rVB	872294	1449198	71.20%	4.223%
94	22.776	3279	3283	3289	rVB7	51375	92806	4.56%	0.270%
95	23.181	3346	3352	3361	rVB	399561	738028	36.26%	2.150%
96	24.368	3546	3554	3563	rVB	428855	1119009	54.98%	3.261%
97	25.249	3694	3704	3711	rBV2	339919	1063936	52.27%	3.100%
98	25.326	3712	3717	3727	rVB2	222424	572202	28.11%	1.667%
99	25.490	3737	3745	3756	rVB2	190366	567863	27.90%	1.655%
100	29.614	4434	4447	4461	rVB	313130	1457355	71.60%	4.246%

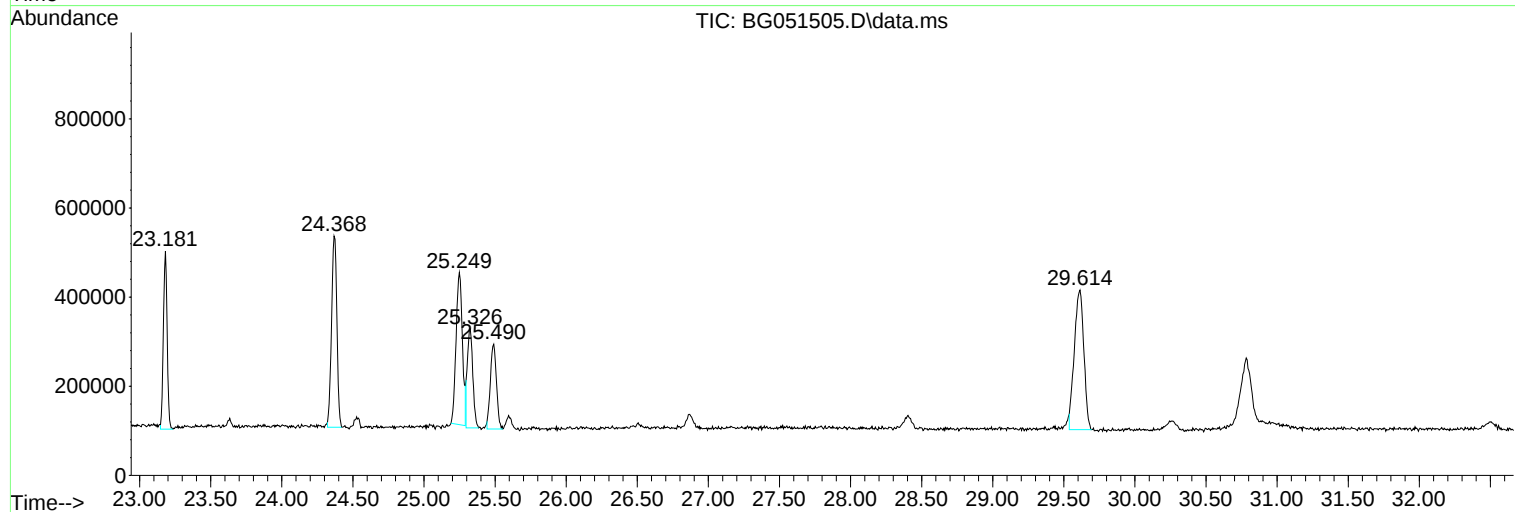
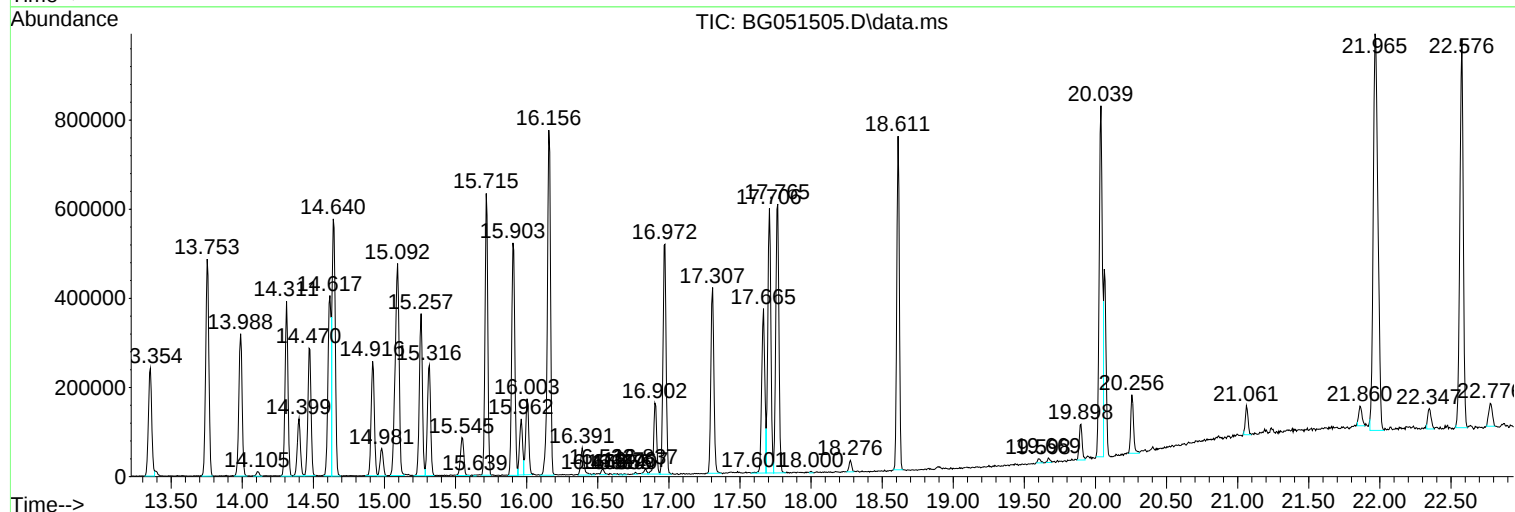
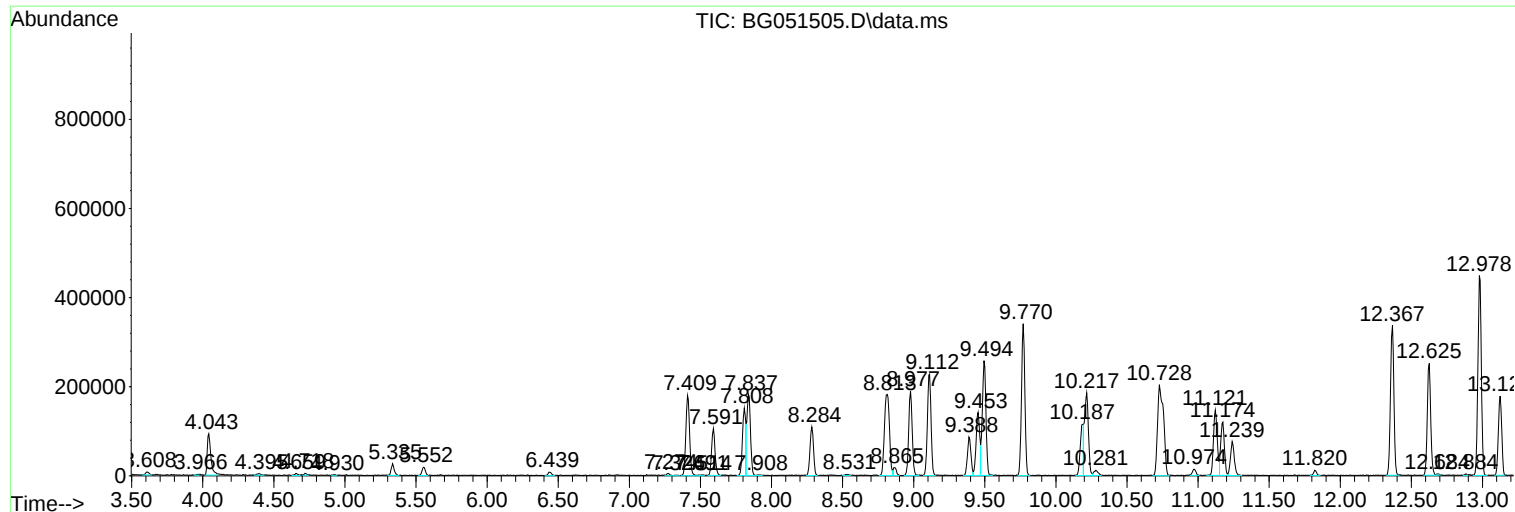
Sum of corrected areas: 34319823

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Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBLS7

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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ALS Vial : 4 Sample Multiplier: 1

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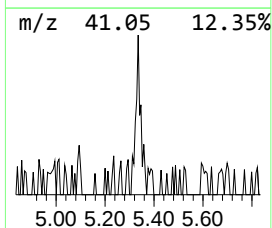
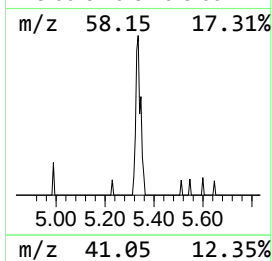
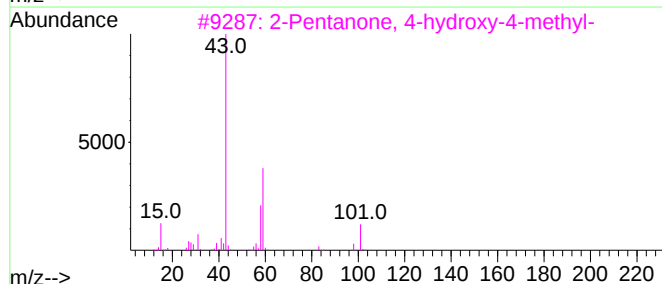
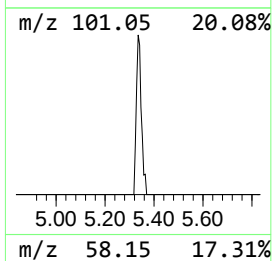
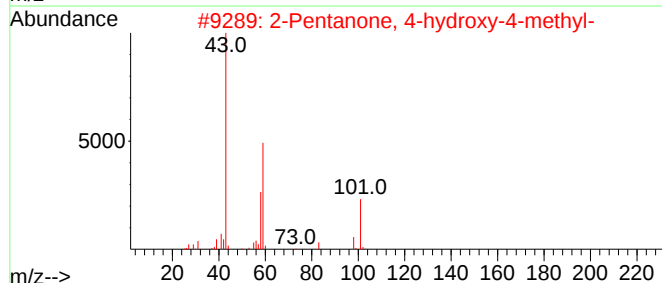
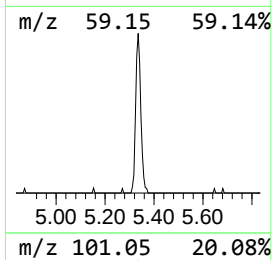
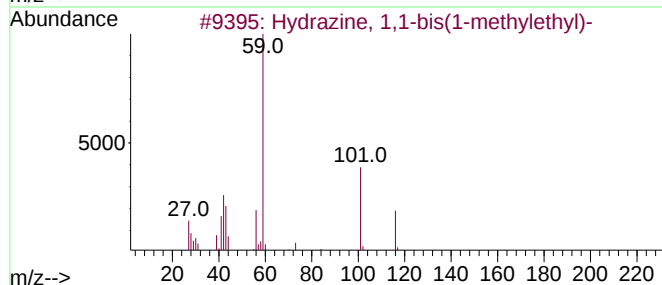
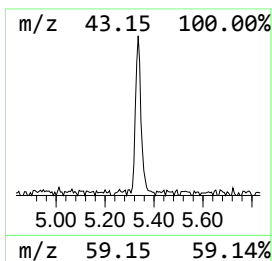
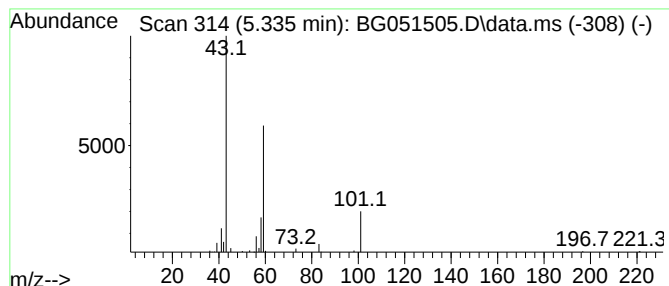
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 unknown-01 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.335	4.19 ng/ul	39166	1,4-Dichlorobenzene-d4	8.284

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hydrazine, 1,1-bis(1-methylethyl)-	116	C6H16N2	000921-14-2	47
2			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45
3			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	38
4			Succinic acid, 2-methoxyethyl te...	372	C21H40O5	1000325-81-3	33
5			3-Octanol	130	C8H18O	000589-98-0	28



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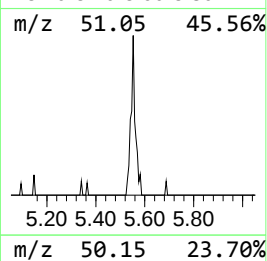
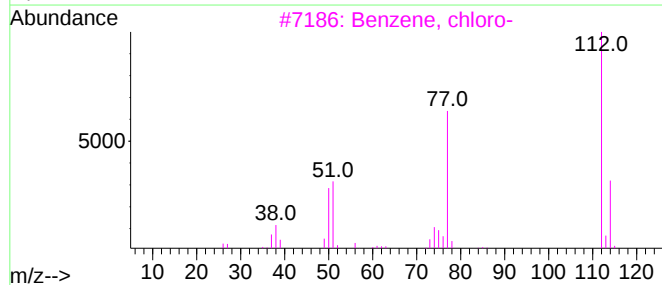
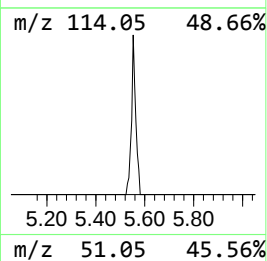
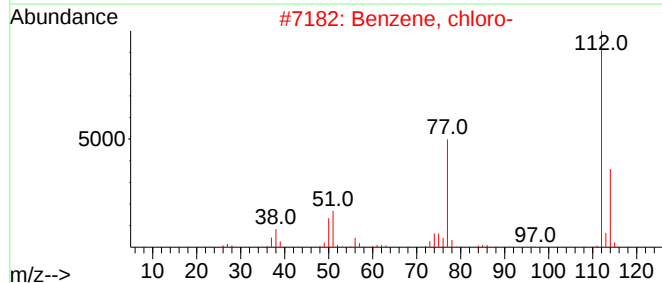
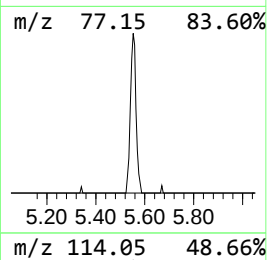
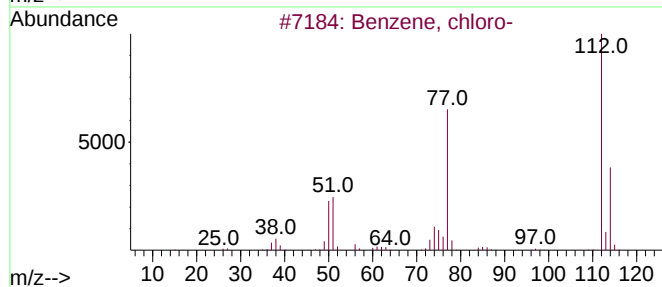
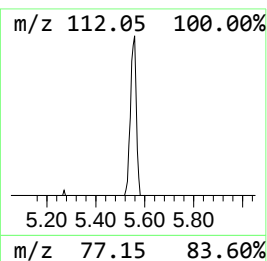
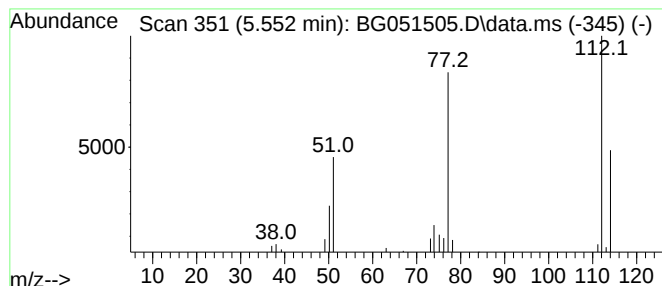
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzene, chloro- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.552	3.45 ng/ul	32218	1,4-Dichlorobenzene-d4	8.284

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, chloro-	112	C6H5Cl	000108-90-7	90
2			Benzene, chloro-	112	C6H5Cl	000108-90-7	86
3			Benzene, chloro-	112	C6H5Cl	000108-90-7	74
4			Benzene, chloro-	112	C6H5Cl	000108-90-7	74
5			Benzene, chloro-	112	C6H5Cl	000108-90-7	64



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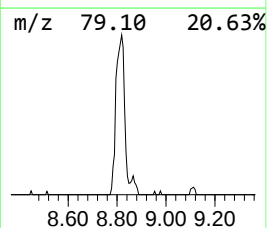
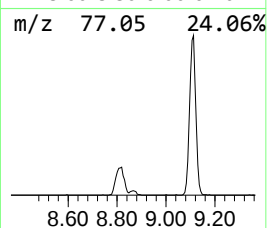
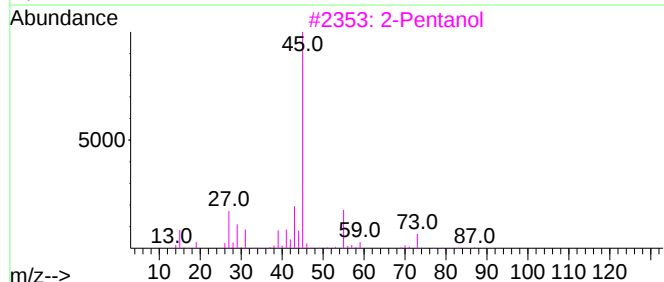
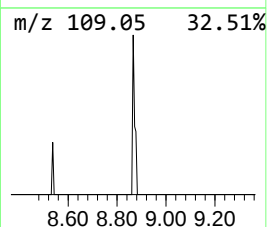
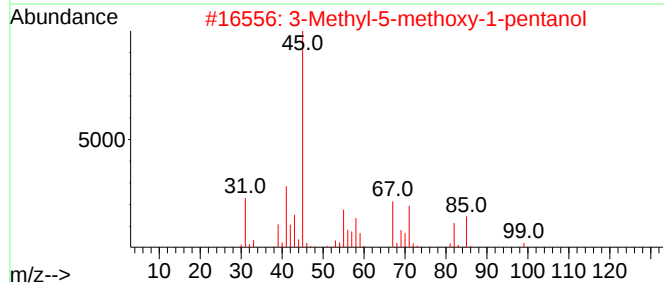
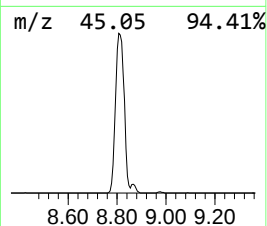
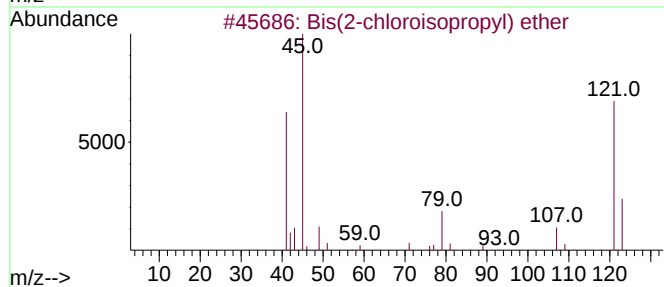
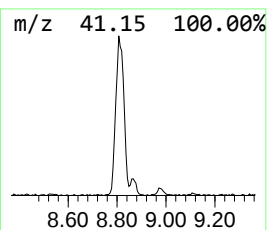
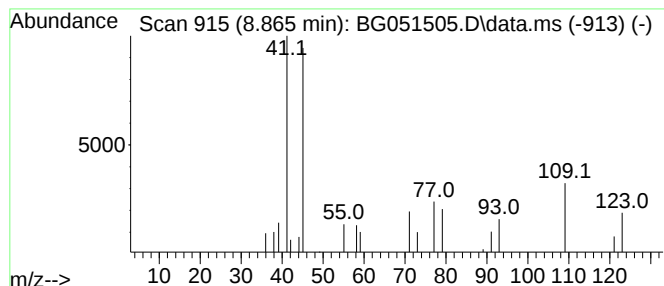
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 unknown-02 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.865	2.70 ng/ul	25196	1,4-Dichlorobenzene-d4	8.284

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Bis(2-chloroisopropyl) ether	170	C6H12Cl2O	039638-32-9	25
2			3-Methyl-5-methoxy-1-pentanol	132	C7H16O2	070928-41-5	9
3			2-Pentanol	88	C5H12O	006032-29-7	9
4			Butane, 1,2,4-trimethoxy-	148	C7H16O3	020637-48-3	9
5			1-Methoxy-5-trifluoroacetoxyhexane	228	C9H15F3O3	1000216-63-2	9



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121421\
Data File : BG051505.D
Acq On : 14 Dec 2021 15:46
Operator : CG/JU
Sample : M5013-05
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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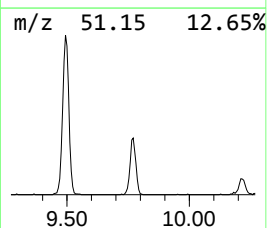
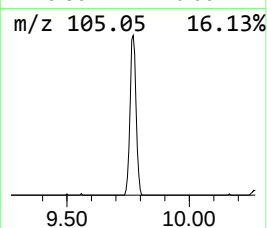
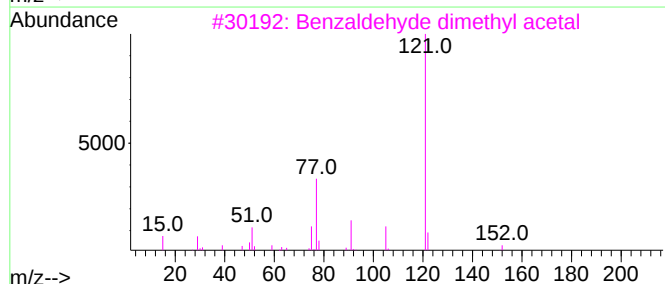
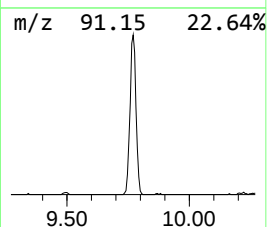
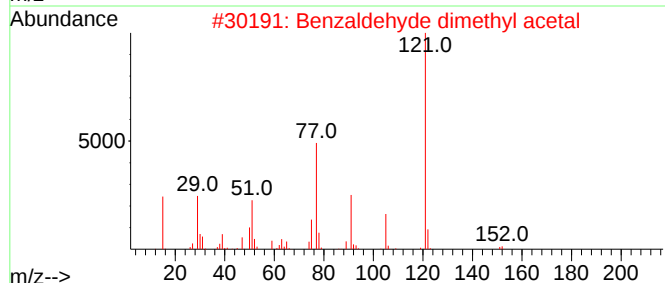
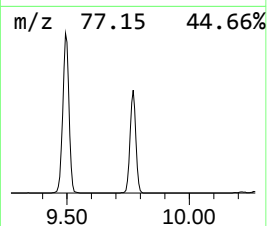
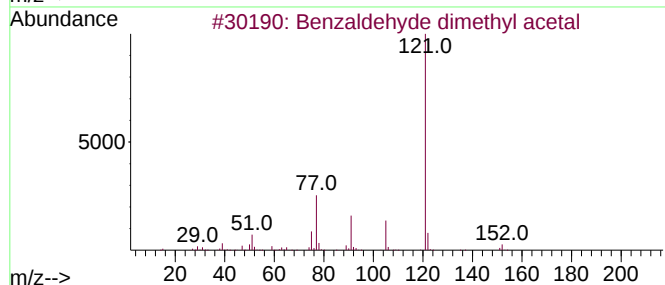
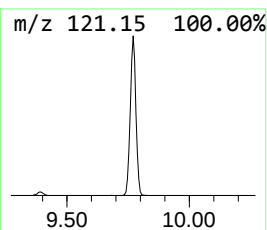
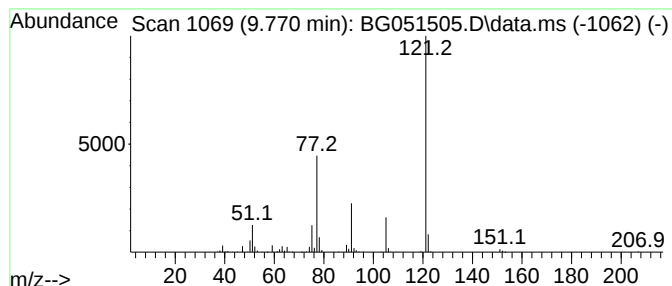
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Benzaldehyde dimethyl acetal Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.770	39.83 ng/ul	555258	Naphthalene-d8	11.121

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzaldehyde dimethyl acetal	152	C9H12O2	001125-88-8	95
2			Benzaldehyde dimethyl acetal	152	C9H12O2	001125-88-8	91
3			Benzaldehyde dimethyl acetal	152	C9H12O2	001125-88-8	90
4			Benzeneacetic acid, .alpha.-meth...	180	C10H12O3	056143-21-6	78
5			Benzeneacetic acid, .alpha.-meth...	180	C10H12O3	056143-21-6	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121421\
Data File : BG051505.D
Acq On : 14 Dec 2021 15:46
Operator : CG/JU
Sample : M5013-05
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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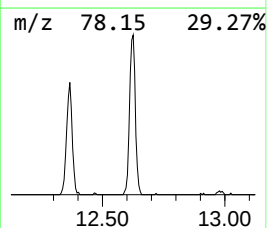
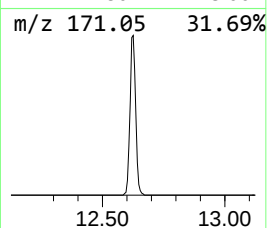
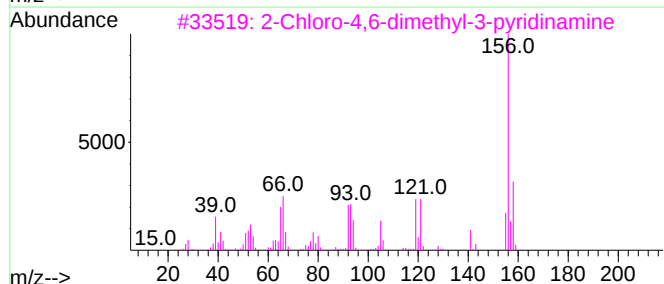
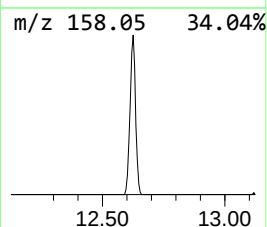
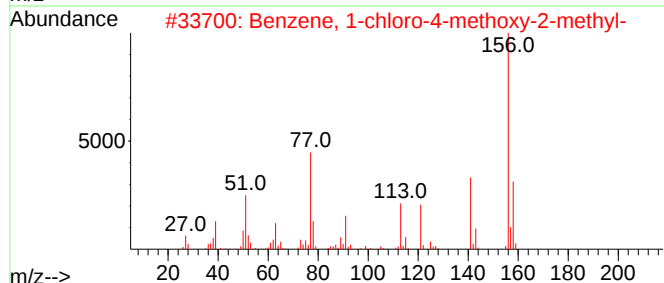
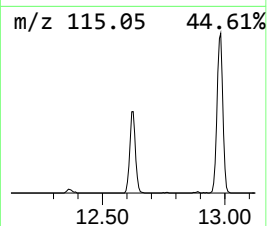
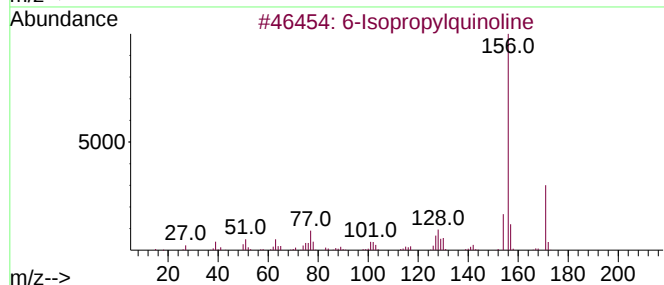
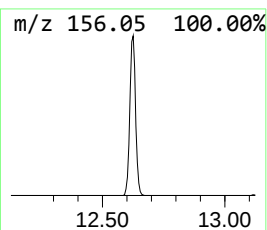
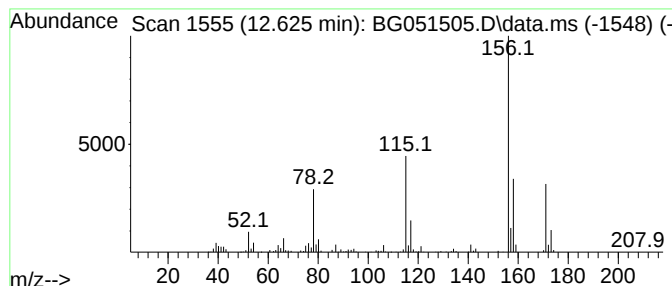
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 unknown-03 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.625	29.87 ng/ul	416437	Naphthalene-d8	11.121

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			6-Isopropylquinoline	171	C12H13N	000135-79-5	43
2			Benzene, 1-chloro-4-methoxy-2-me...	156	C8H9ClO	013334-71-9	38
3			2-Chloro-4,6-dimethyl-3-pyridina...	156	C7H9ClN2	140413-40-7	38
4			Chloroxylenol	156	C8H9ClO	000088-04-0	32
5			Tebuthiuron	228	C9H16N4O5	034014-18-1	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121421\
Data File : BG051505.D
Acq On : 14 Dec 2021 15:46
Operator : CG/JU
Sample : M5013-05
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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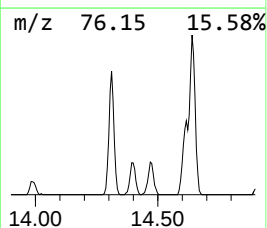
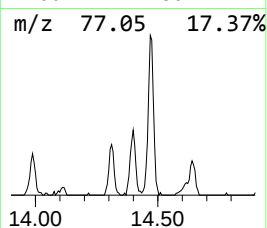
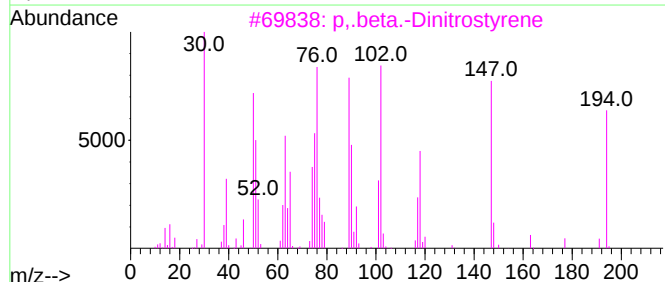
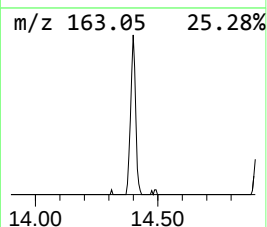
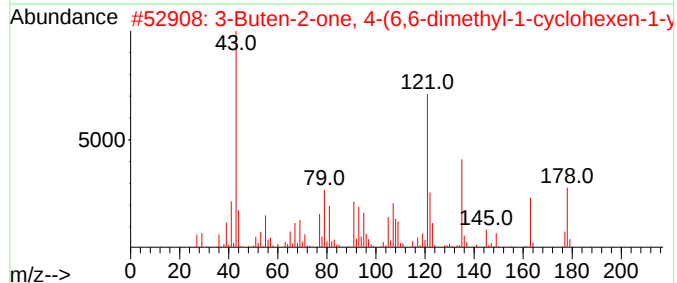
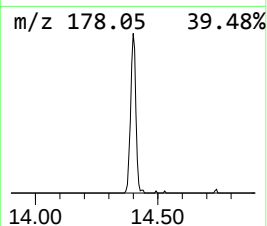
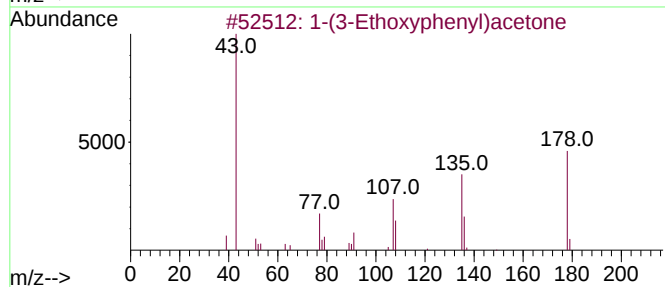
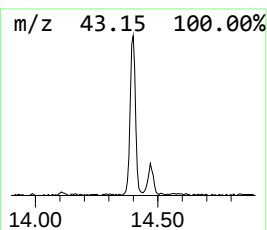
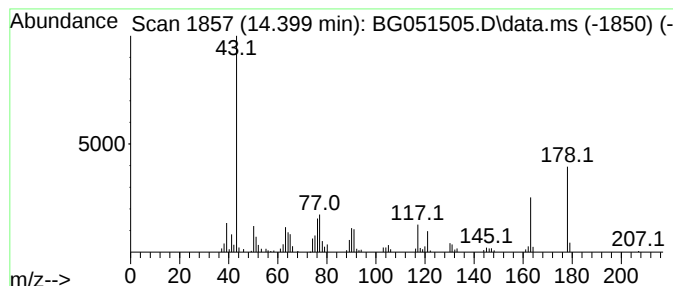
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 unknown-04 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.399	9.81 ng/ul	198026	Acenaphthene-d10	14.916

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-(3-Ethoxyphenyl)acetone	178	C11H14O2	023037-47-0	25
2			3-Buten-2-one, 4-(6,6-dimethyl-1...	178	C12H18O	065133-79-1	23
3			p,.beta.-Dinitrostyrene	194	C8H6N2O4	003156-41-0	10
4			2,6-Diacetylpyridine	163	C9H9NO2	001129-30-2	9
5			2-Bromophenethyl alcohol, n-prop...	242	C11H15BrO	1000394-87-6	9



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121421\
Data File : BG051505.D
Acq On : 14 Dec 2021 15:46
Operator : CG/JU
Sample : M5013-05
Misc :
ALS Vial : 4 Sample Multiplier: 1

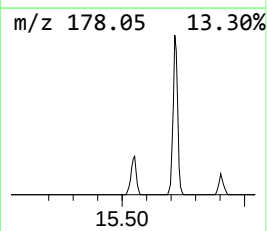
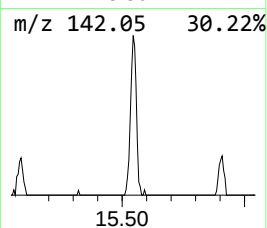
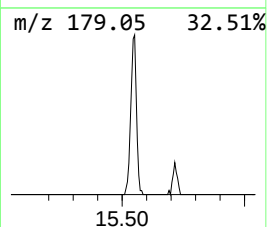
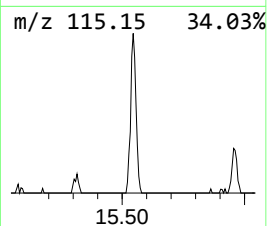
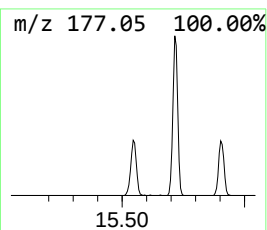
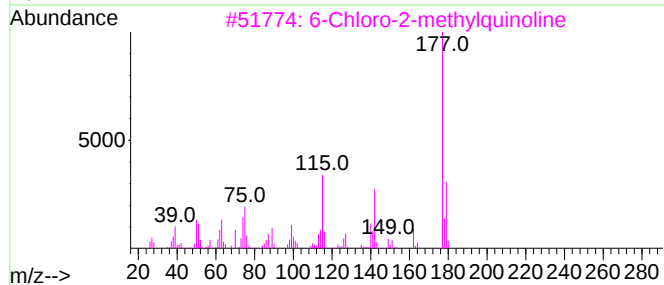
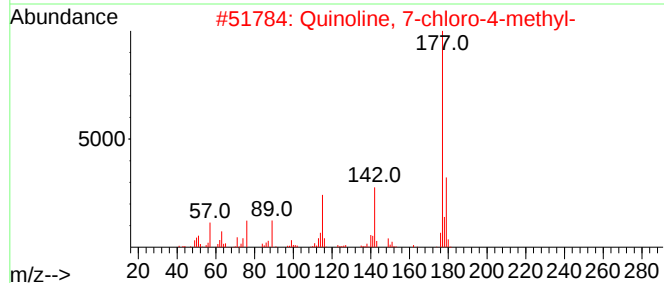
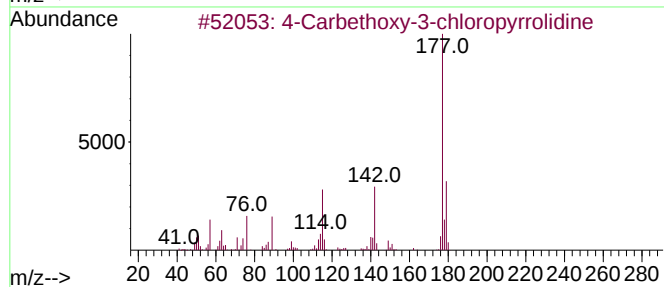
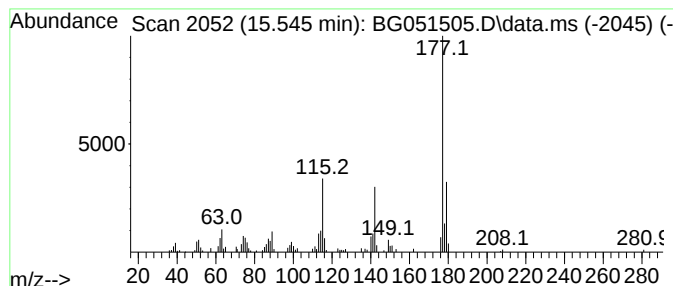
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 4-Carboxy-3-chloropyrrol... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD		R.T.	
15.545	6.93 ng/ul	139952	Acenaphthene-d10		14.916	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	4-Carbethoxy-3-chloropyrrolidine		177	C7H12ClNO2	1010255-97-9	93
2	Quinoline, 7-chloro-4-methyl-		177	C10H8ClN	040941-53-5	93
3	6-Chloro-2-methylquinoline		177	C10H8ClN	000092-46-6	91
4	Quinoline, 2-chloro-4-methyl-		177	C10H8ClN	000634-47-9	91
5	2-Chloro-3-methylquinoline		177	C10H8ClN	057876-69-4	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121421\
Data File : BG051505.D
Acq On : 14 Dec 2021 15:46
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Sample : M5013-05
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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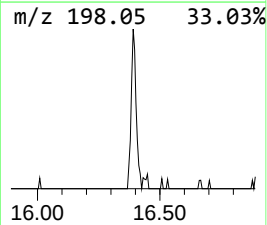
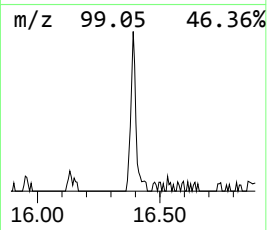
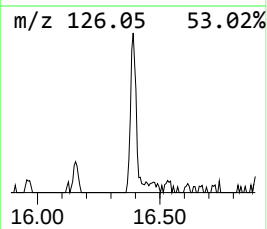
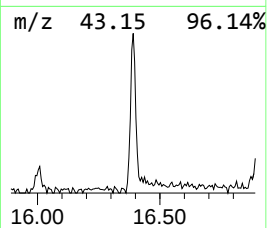
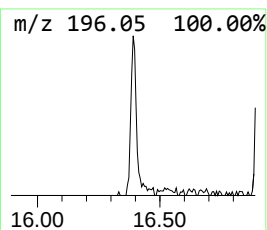
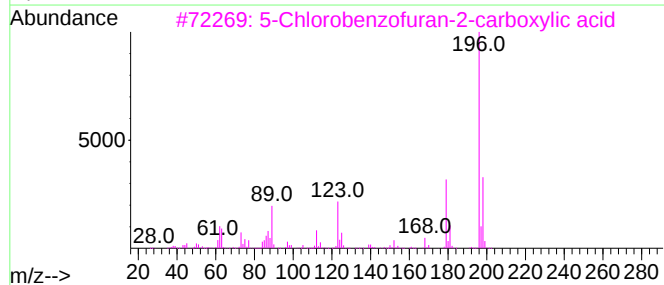
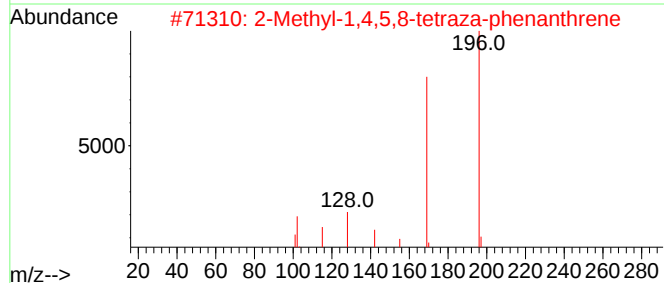
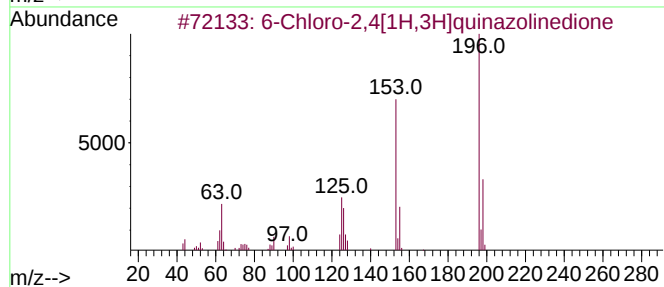
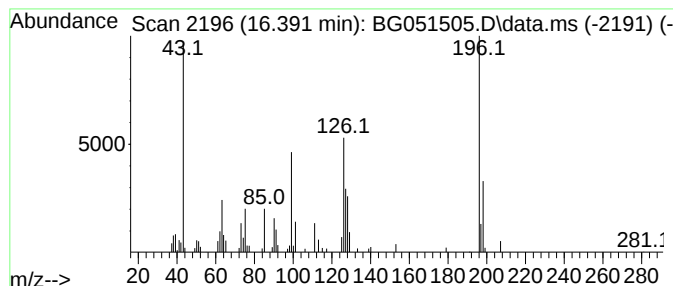
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 unknown-05 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.391	3.47 ng/ul	98749	Phenanthrene-d10	17.665

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			6-Chloro-2,4[1H,3H]quinazolinedione	196	C8H5ClN2O2	001640-60-4	49
2			2-Methyl-1,4,5,8-tetraza-phenant...	196	C11H8N4	103538-90-5	43
3			5-Chlorobenzofuran-2-carboxylic ...	196	C9H5ClO3	1000484-31-1	38
4			4-Amino-2-bromobenzonitrile	196	C7H5BrN2	053312-82-6	38
5			2-Chloro-4-methyl-6-(trifluorome...	196	C6H4ClF3N2	241164-09-0	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121421\
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Sample : M5013-05
Misc :
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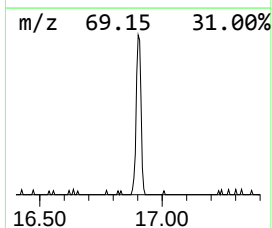
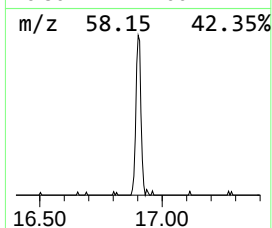
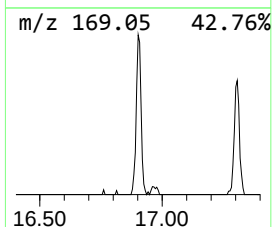
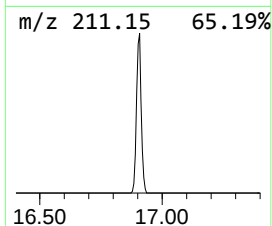
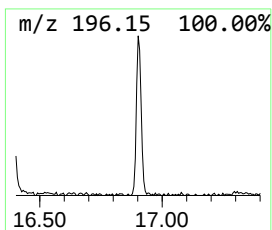
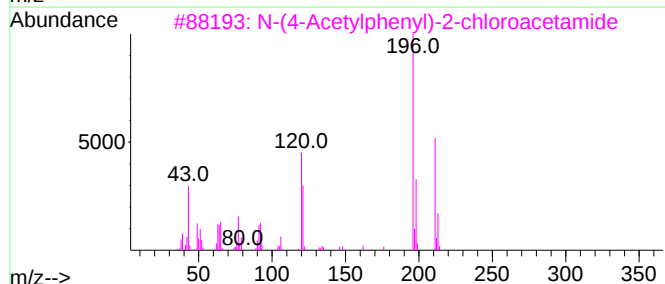
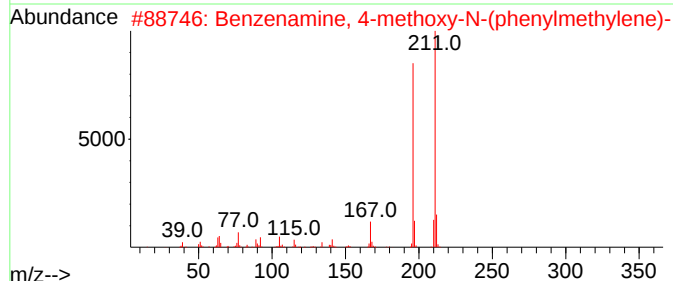
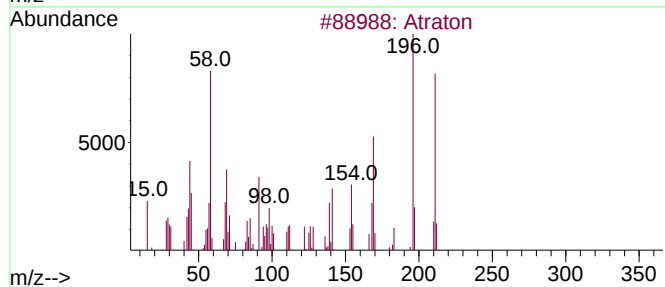
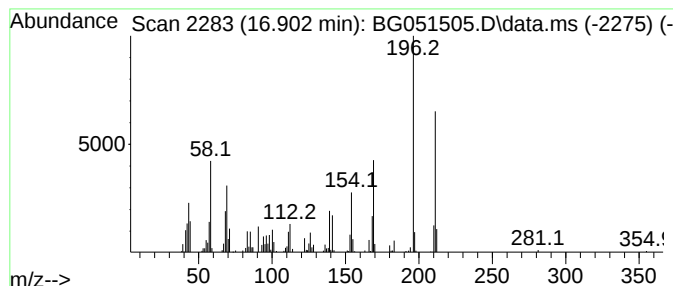
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Atraton Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.902	8.12 ng/ul	231058	Phenanthrene-d10	17.665

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Atraton			211	C9H17N5O	001610-17-9	91
2	Benzenamine, 4-methoxy-N-(phenyl...			211	C14H13NO	000783-08-4	46
3	N-(4-Acetylphenyl)-2-chloroaceta...			211	C10H10ClNO2	038283-38-4	43
4	2-Acetyl-4-methoxy-6-nitrophenol			211	C9H9NO5	090564-25-3	38
5	Benzenamine, 4-methoxy-N-(phenyl...			211	C14H13NO	000783-08-4	30



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121421\
Data File : BG051505.D
Acq On : 14 Dec 2021 15:46
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Misc :
ALS Vial : 4 Sample Multiplier: 1

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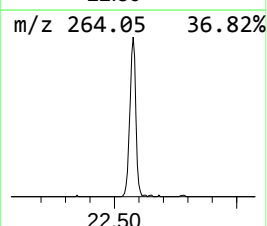
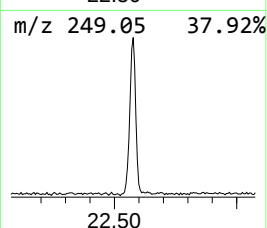
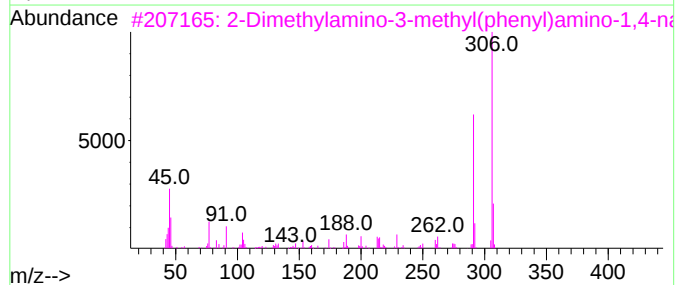
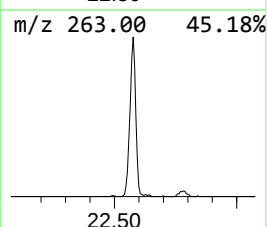
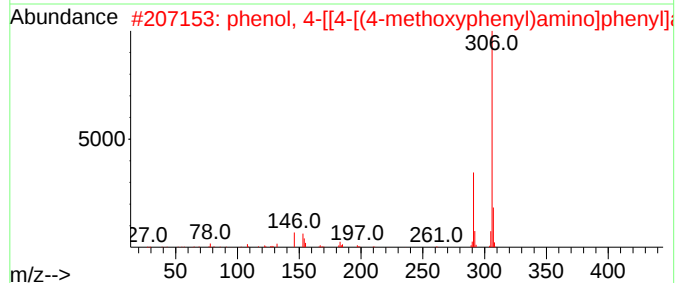
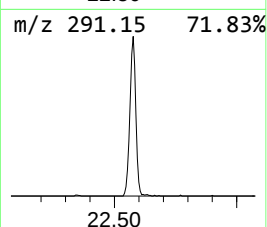
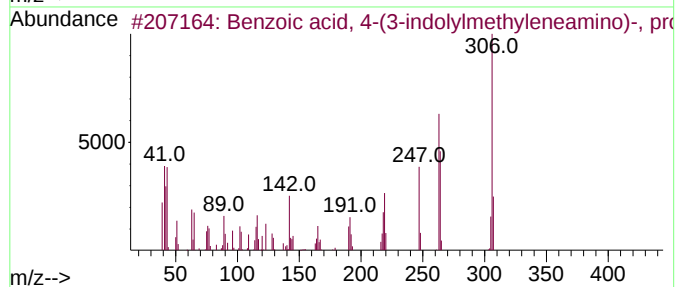
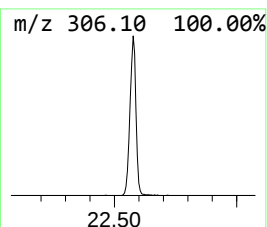
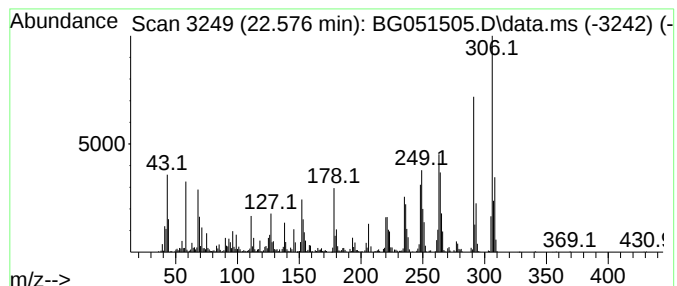
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 10 unknown-06 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.576	14.24 ng/ul	1449200	Chrysene-d12	21.983

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzoic acid, 4-(3-indolylmethyl...	306	C19H18N2O2	304669-02-1	46
2			phenol, 4-[[4-[(4-methoxyphenyl)...	306	C19H18N2O2	1000397-62-8	46
3			2-Dimethylamino-3-methyl(phenyl)...	306	C19H18N2O2	134614-10-1	45
4			Pyrido[2,3-b]quinolin-5-one, 7-c...	306	C18H11ClN2O	069750-07-8	42
5			Dibenzo[def,mno]chrysene-6,12-dione	306	C22H10O2	000641-13-4	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121421\
Data File : BG051505.D
Acq On : 14 Dec 2021 15:46
Operator : CG/JU
Sample : M5013-05
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBLS7

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
unknown-01	5.335	4.2	ng/ul	39166	1	8.284	186830	20.0
Benzene, chloro-	5.552	3.5	ng/ul	32218	1	8.284	186830	20.0
unknown-02	8.865	2.7	ng/ul	25196	1	8.284	186830	20.0
Benzaldehyde di...	9.770	39.8	ng/ul	555258	2	11.121	278826	20.0
unknown-03	12.625	29.9	ng/ul	416437	2	11.121	278826	20.0
unknown-04	14.399	9.8	ng/ul	198026	3	14.916	403613	20.0
4-Carbethoxy-3-...	15.545	6.9	ng/ul	139952	3	14.916	403613	20.0
unknown-05	16.391	3.5	ng/ul	98749	4	17.665	569146	20.0
Atraton	16.902	8.1	ng/ul	231058	4	17.665	569146	20.0
unknown-06	22.576	14.2	ng/ul	1449200	5	21.983	2035350	20.0