

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG053122\
 Data File : BG053776.D
 Acq On : 1 Jun 2022 2:35
 Operator : CG/JU
 Sample : N3039-11
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0JR4

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-BG053122.M
 Title : SVOA CALIBRATION

Signal : TIC: BG053776.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.169	24	31	43	rBV	1867988	4191259	10.59%	2.743%
2	3.269	43	48	65	rVB	411333	822795	2.08%	0.539%
3	5.473	404	423	428	rBV3	8868878	37938923	95.83%	24.833%
4	6.783	640	646	655	rVB2	22217	43578	0.11%	0.029%
5	7.247	720	725	737	rVB3	22160	49037	0.12%	0.032%
6	7.435	748	757	767	rVB	68770	144268	0.36%	0.094%
7	7.635	782	791	803	rVB	72472	157187	0.40%	0.103%
8	8.017	844	856	867	rBV	2322182	6522231	16.47%	4.269%
9	8.210	867	889	893	rBV	8752721	33152079	83.74%	21.700%
10	8.557	923	948	952	rBV	8991983	39591461	100.00%	25.915%
11	8.798	979	989	997	rVB	63540	116154	0.29%	0.076%
12	9.262	1061	1068	1071	rBV	96392	180767	0.46%	0.118%
13	9.309	1071	1076	1084	rVB	240299	447185	1.13%	0.293%
14	9.603	1120	1126	1133	rVB	79766	140310	0.35%	0.092%
15	9.920	1175	1180	1186	rVV	42004	85762	0.22%	0.056%
16	9.985	1186	1191	1199	rVV	68191	133068	0.34%	0.087%
17	10.525	1275	1283	1291	rBV	112028	213085	0.54%	0.139%
18	10.819	1317	1333	1338	rBV	5970814	16393820	41.41%	10.731%
19	10.931	1345	1352	1357	rBV	96508	164751	0.42%	0.108%
20	11.042	1363	1371	1380	rVB	106674	206841	0.52%	0.135%
21	11.313	1404	1417	1426	rVB	2251514	4438956	11.21%	2.906%
22	11.501	1442	1449	1458	rVB	132681	240926	0.61%	0.158%
23	11.712	1477	1485	1493	rVB2	32109	58277	0.15%	0.038%
24	12.923	1676	1691	1698	rVB	275228	538705	1.36%	0.353%
25	13.528	1786	1794	1802	rBV	766801	1257701	3.18%	0.823%
26	14.121	1885	1895	1902	rBV	234078	379337	0.96%	0.248%
27	14.415	1938	1945	1953	rVB	260437	427112	1.08%	0.280%
28	14.721	1990	1997	2003	rVB	145586	238000	0.60%	0.156%
29	15.713	2158	2166	2174	rBV	347972	582834	1.47%	0.381%
30	15.813	2177	2183	2191	rBV	79251	122062	0.31%	0.080%
31	17.464	2456	2464	2471	rVB	197520	304483	0.77%	0.199%
32	17.564	2474	2481	2491	rVB	397802	633608	1.60%	0.415%
33	18.827	2687	2696	2703	rBV2	143675	208265	0.53%	0.136%
34	19.844	2863	2869	2877	rBV	567056	836762	2.11%	0.548%
35	21.389	3128	3132	3139	rBV3	26396	44311	0.11%	0.029%
36	21.742	3186	3192	3200	rBV2	245945	410800	1.04%	0.269%

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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-BG053122.M
Title : SVOA CALIBRATION

37	24.791	3700	3711	3721	rBV2	318993	909972	2.30%	0.596%
38	25.014	3740	3749	3761	rVB2	134253	394558	1.00%	0.258%
39	25.578	3843	3845	3856	rVB2	23168	53495	0.14%	0.035%

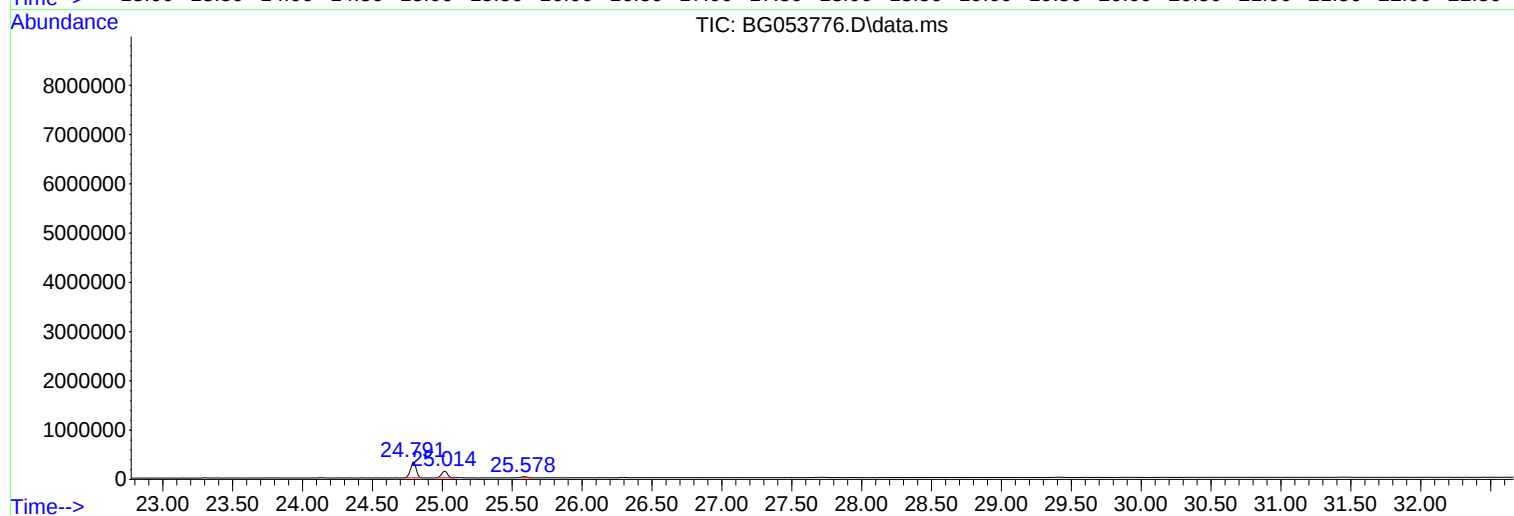
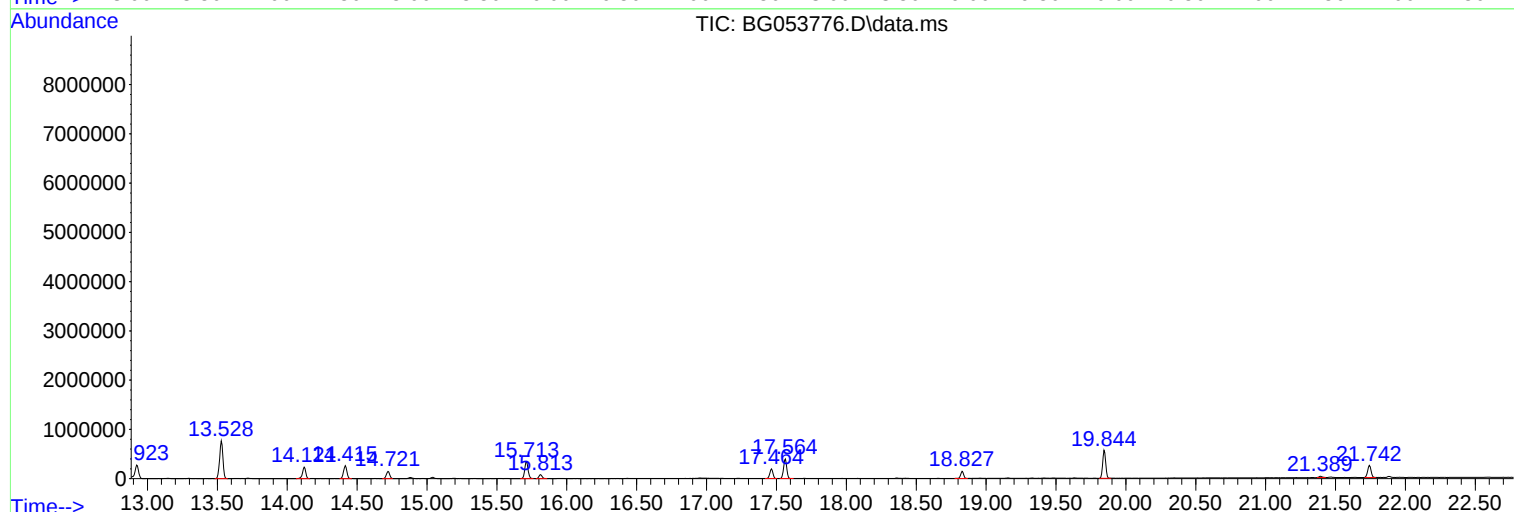
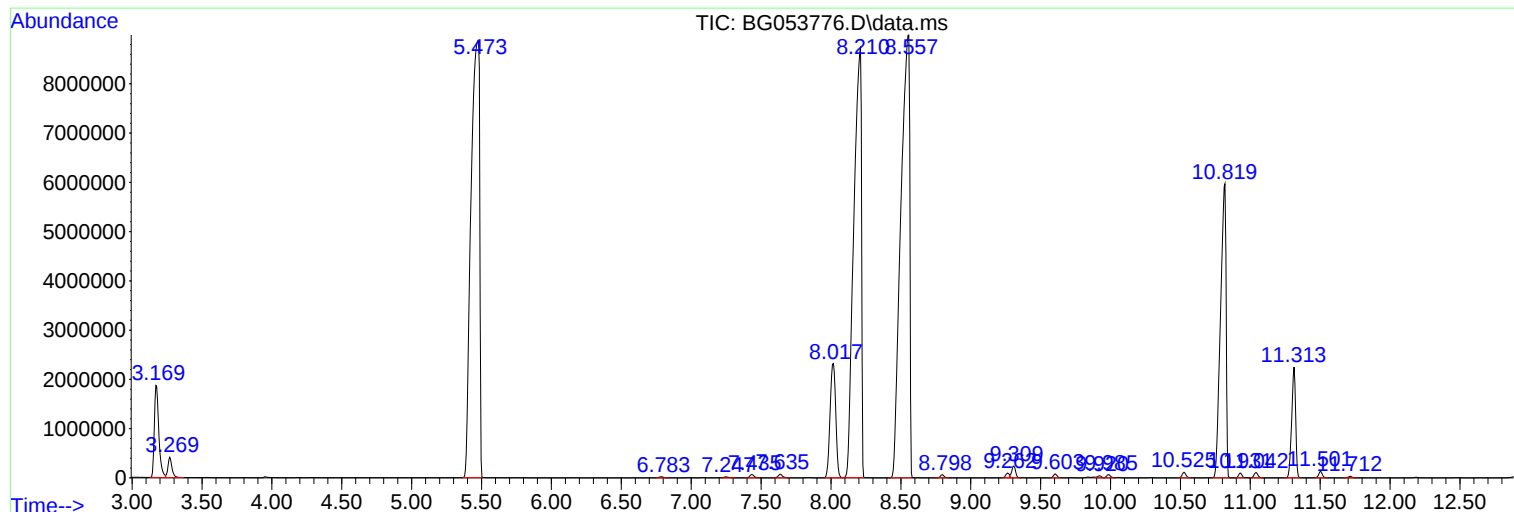
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG053122.M
Quant Title : SVOA CALIBRATION

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



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Operator : CG/JU
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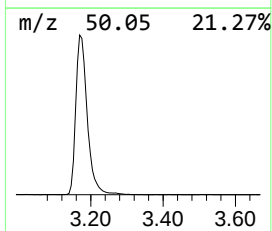
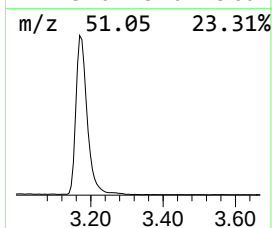
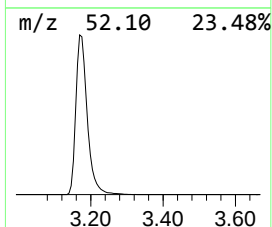
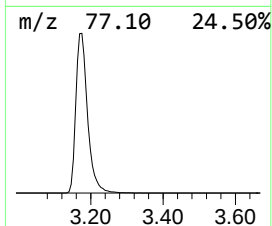
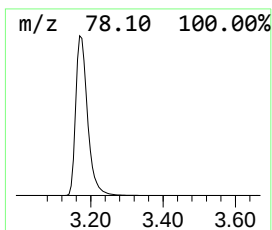
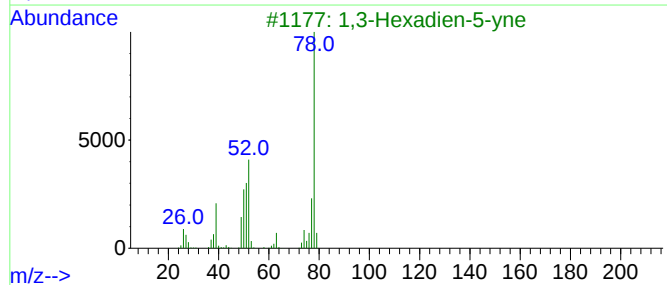
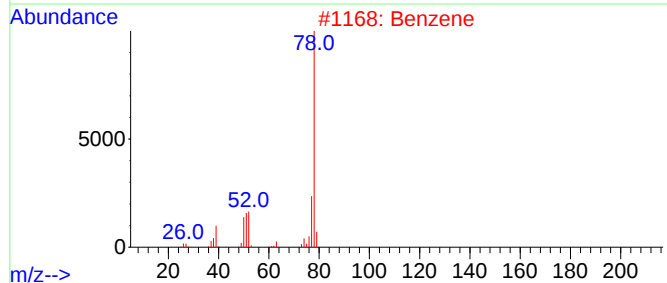
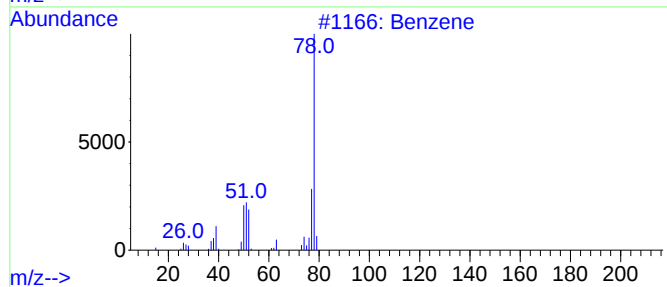
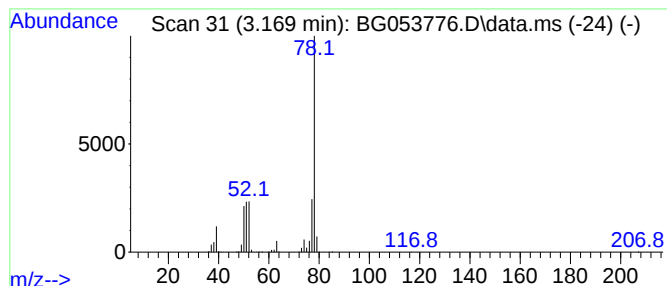
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG053122.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzene Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.169	2.53 ng/ul	4191260	1,4-Dichlorobenzene-d4	8.140

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene	78	C6H6	000071-43-2	95
2			Benzene	78	C6H6	000071-43-2	95
3			1,3-Hexadien-5-yne	78	C6H6	010420-90-3	91
4			Benzene	78	C6H6	000071-43-2	91
5			Benzene	78	C6H6	000071-43-2	91



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ALS Vial : 16 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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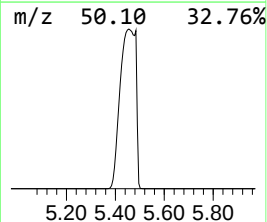
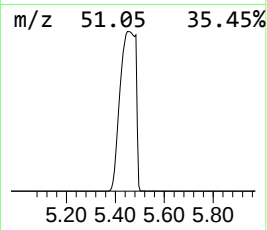
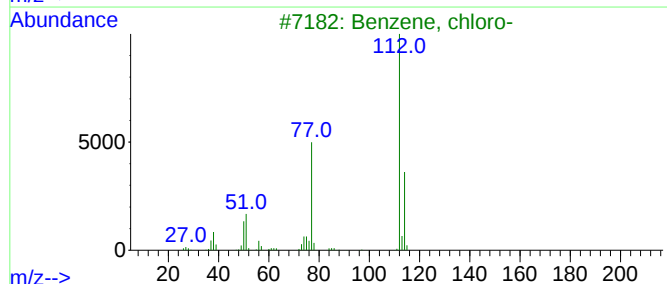
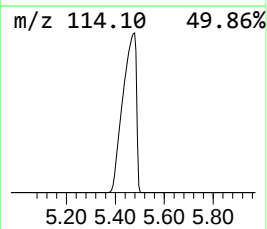
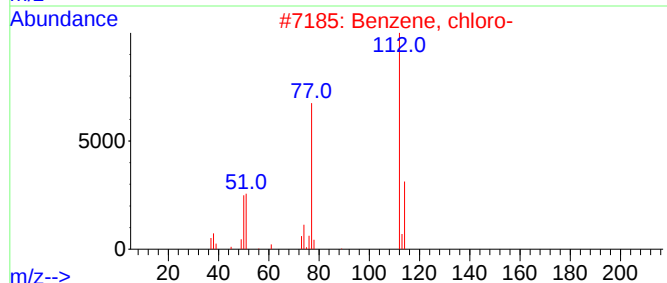
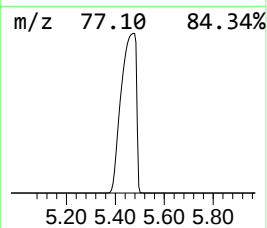
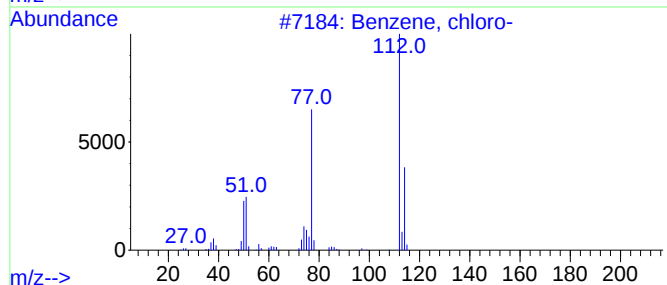
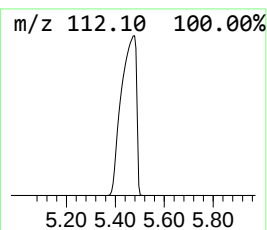
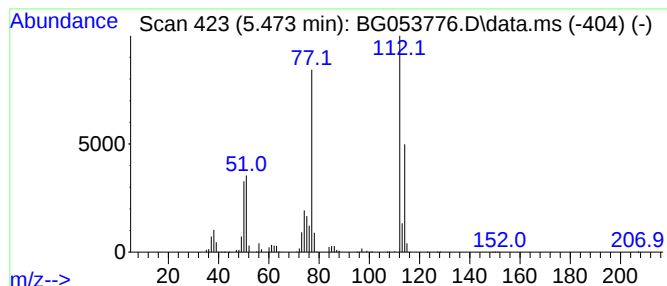
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG053122.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzene, chloro- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.473	22.89 ng/ul	37938900	1,4-Dichlorobenzene-d4	8.140

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, chloro-	112	C6H5Cl	000108-90-7	96
2			Benzene, chloro-	112	C6H5Cl	000108-90-7	94
3			Benzene, chloro-	112	C6H5Cl	000108-90-7	91
4			Benzene, chloro-	112	C6H5Cl	000108-90-7	87
5			Benzene, chloro-	112	C6H5Cl	000108-90-7	87



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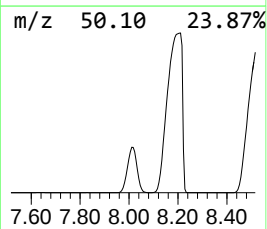
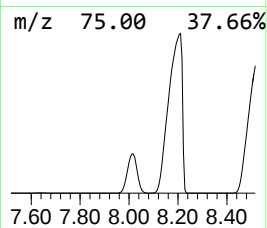
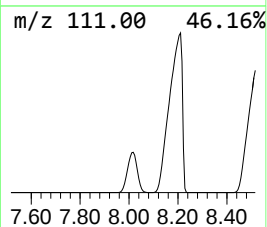
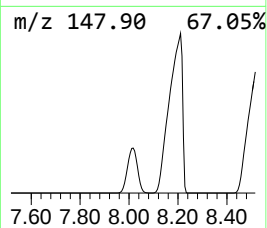
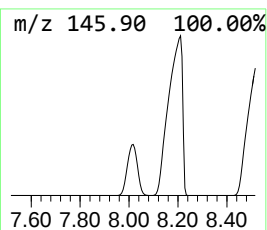
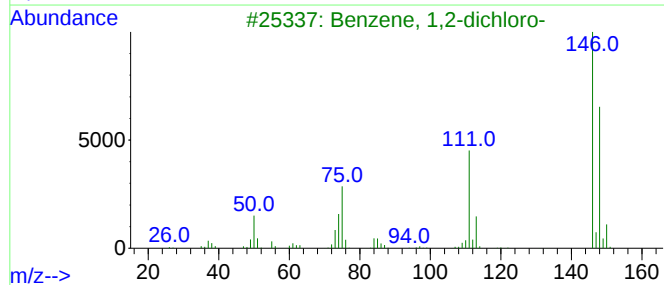
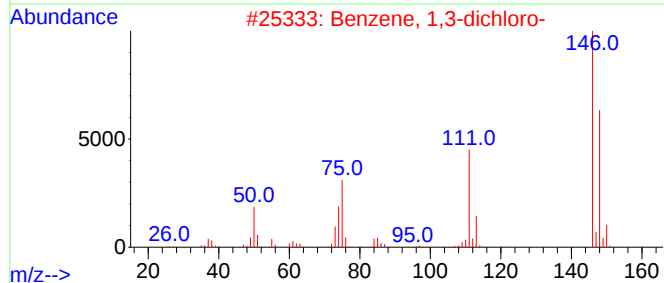
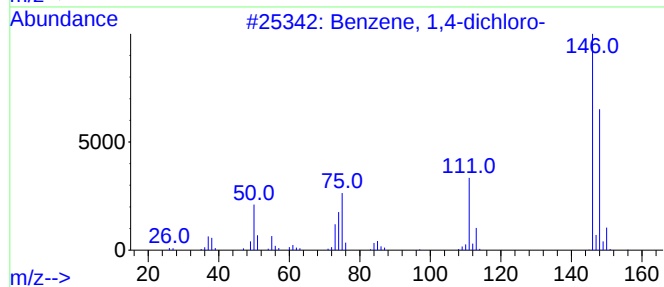
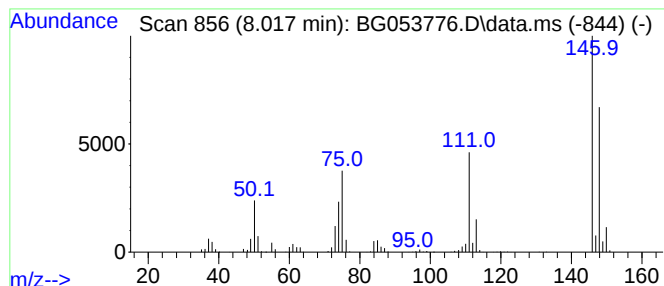
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG053122.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Benzene, 1,4-dichloro- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.017	3.93 ng/ul	6522230	1,4-Dichlorobenzene-d4	8.140

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	98
2			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	98
3			Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	98
4			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	97
5			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	97



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Instrument :
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ClientSampleId :
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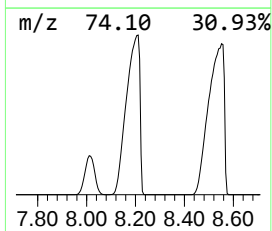
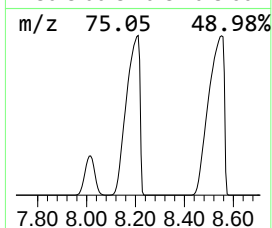
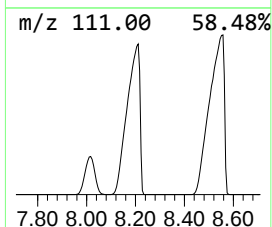
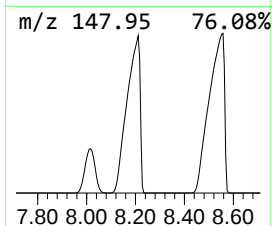
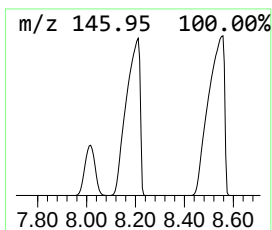
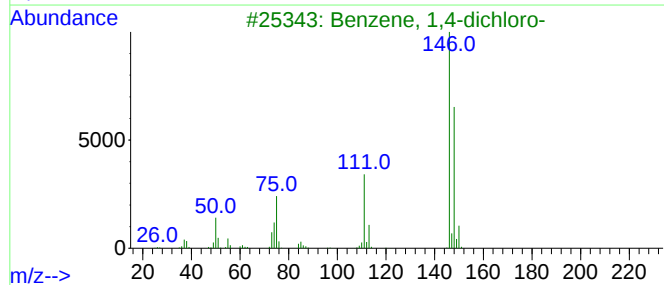
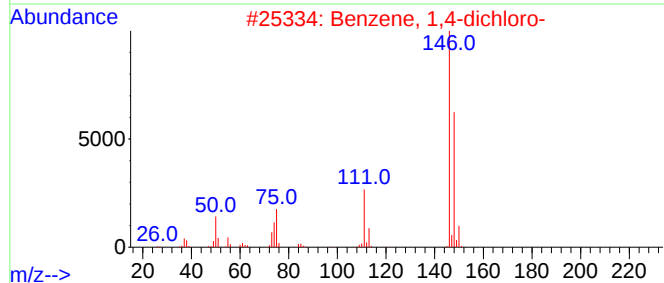
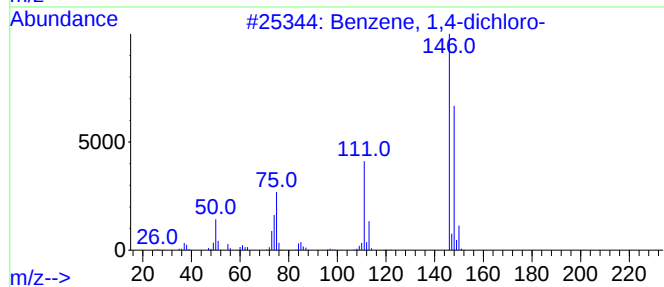
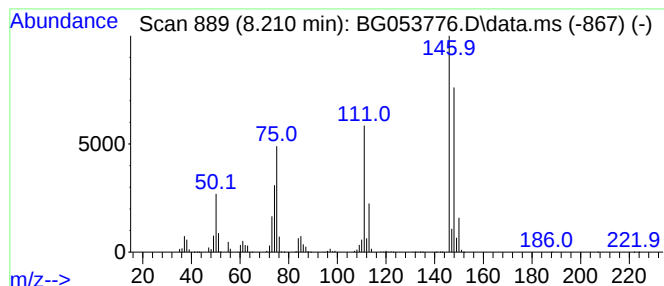
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Benzene, 1,3-dichloro- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.210	20.00 ng/ul	33152100	1,4-Dichlorobenzene-d4	8.140

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	97
2			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	96
3			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	96
4			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	96
5			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	96



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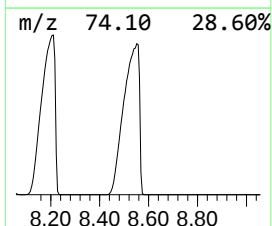
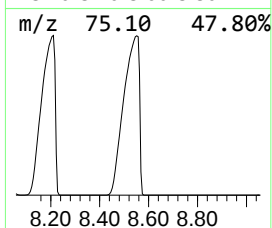
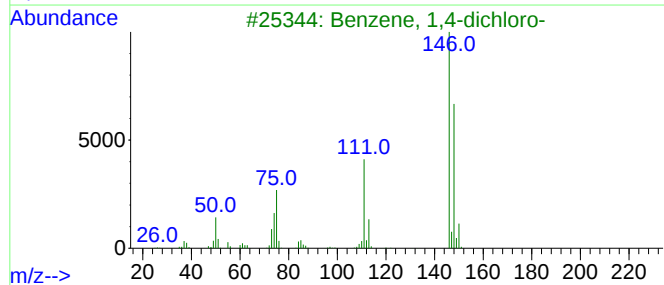
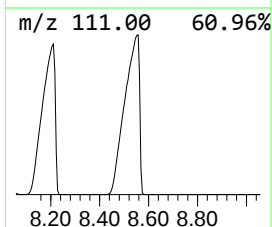
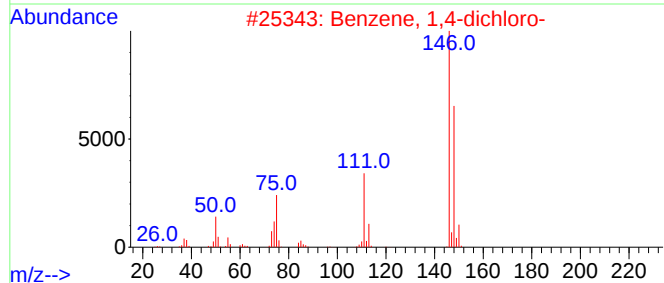
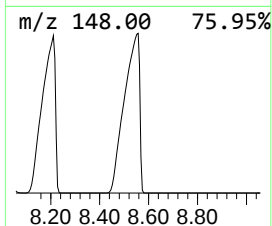
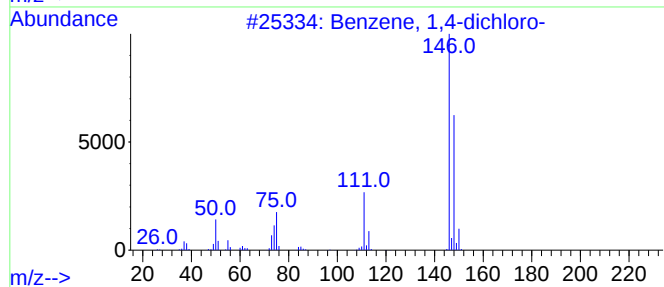
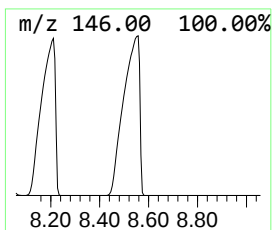
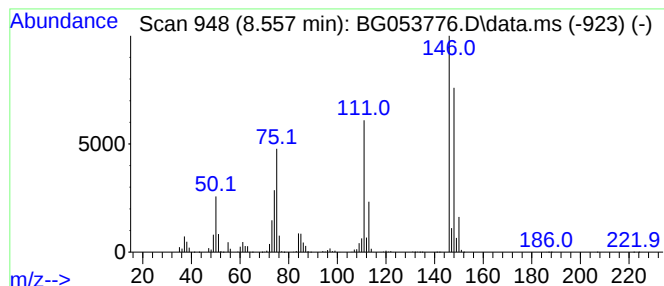
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG053122.M
Quant Title : SVOA CALIBRATION

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

Peak Number 5 Benzene, 1,2-dichloro- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.557	23.88 ng/ul	39591500	1,4-Dichlorobenzene-d4	8.140

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	96
2			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	96
3			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	96
4			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	96
5			Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG053122\
Data File : BG053776.D
Acq On : 1 Jun 2022 2:35
Operator : CG/JU
Sample : N3039-11
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0JR4

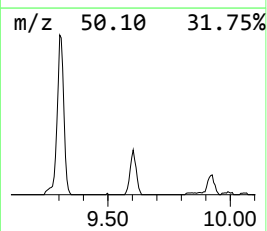
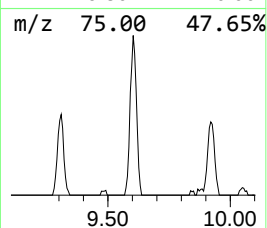
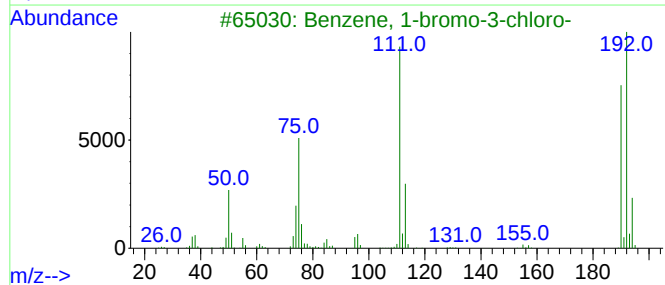
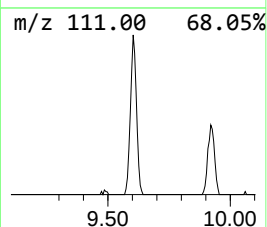
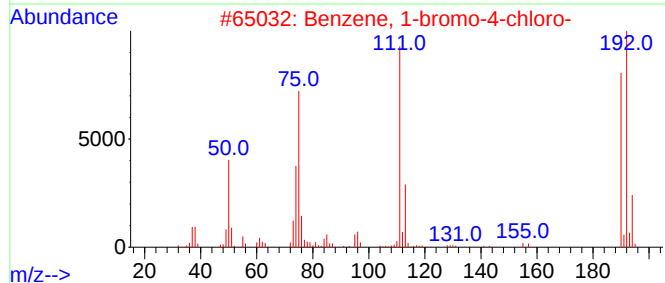
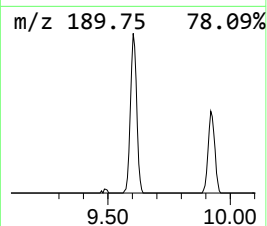
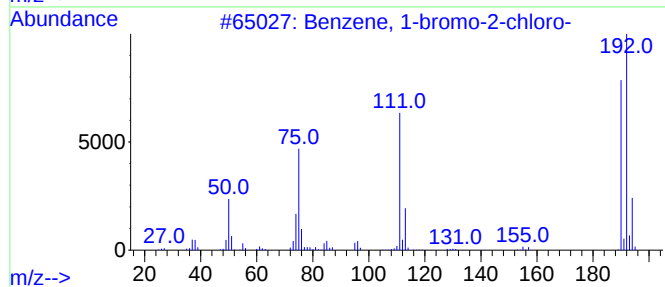
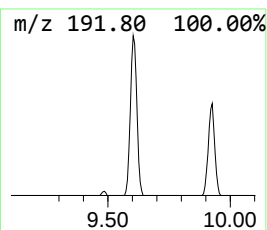
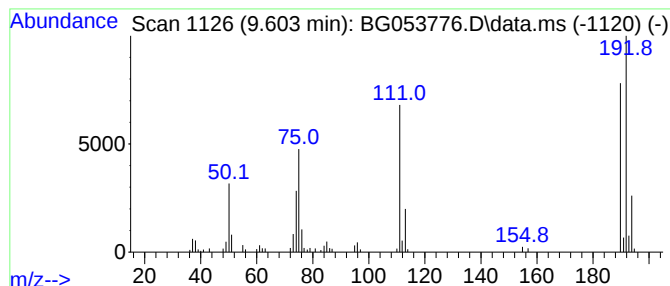
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Benzene, 1-bromo-2-chloro- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.603	17.03 ng/ul	140310	Naphthalene-d8	10.931

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-bromo-2-chloro-	190	C6H4BrCl	000694-80-4	98
2			Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	97
3			Benzene, 1-bromo-3-chloro-	190	C6H4BrCl	000108-37-2	97
4			Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	96
5			Benzene, 1-bromo-3-chloro-	190	C6H4BrCl	000108-37-2	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG053122\
Data File : BG053776.D
Acq On : 1 Jun 2022 2:35
Operator : CG/JU
Sample : N3039-11
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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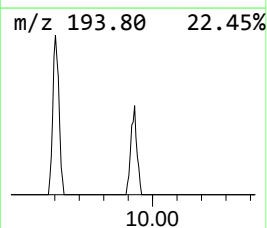
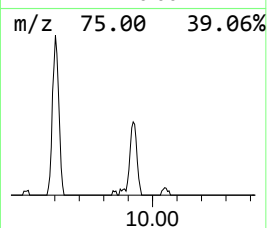
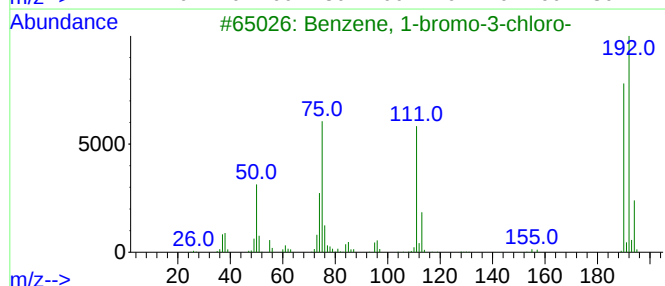
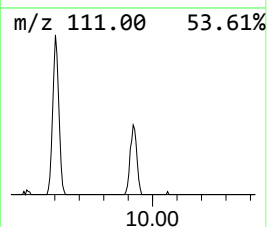
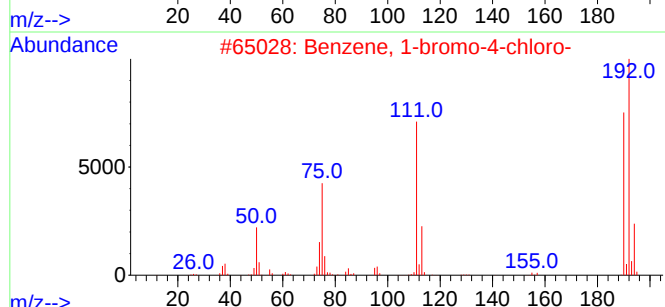
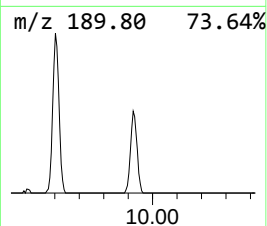
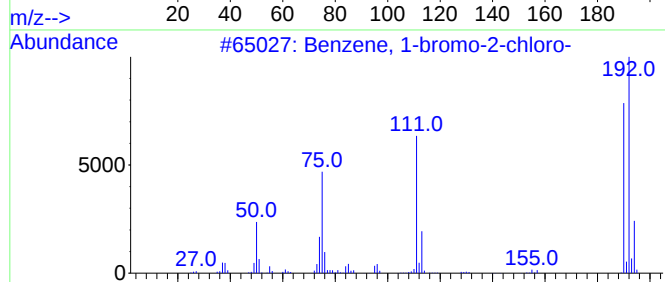
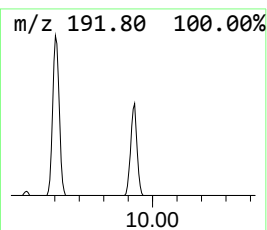
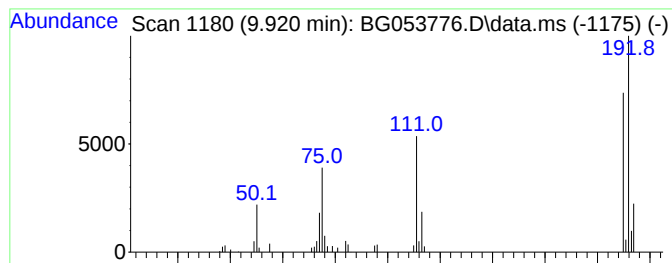
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Benzene, 1-bromo-4-chloro- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.920	10.41 ng/ul	85762	Naphthalene-d8	10.931

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-bromo-2-chloro-	190	C6H4BrCl	000694-80-4	98
2			Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	98
3			Benzene, 1-bromo-3-chloro-	190	C6H4BrCl	000108-37-2	96
4			Benzene, 1-bromo-2-chloro-	190	C6H4BrCl	000694-80-4	96
5			Benzene, 1-bromo-3-chloro-	190	C6H4BrCl	000108-37-2	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG053122\
Data File : BG053776.D
Acq On : 1 Jun 2022 2:35
Operator : CG/JU
Sample : N3039-11
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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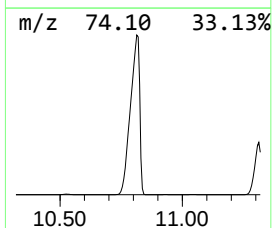
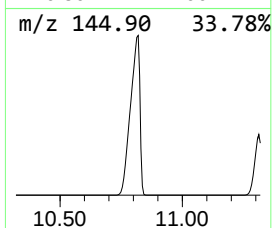
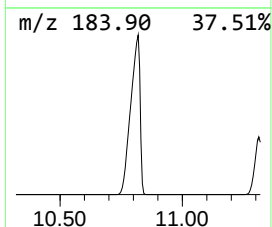
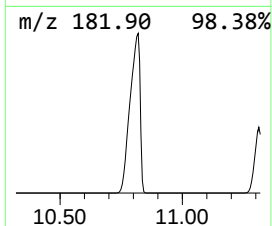
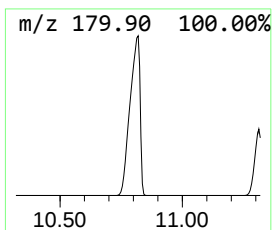
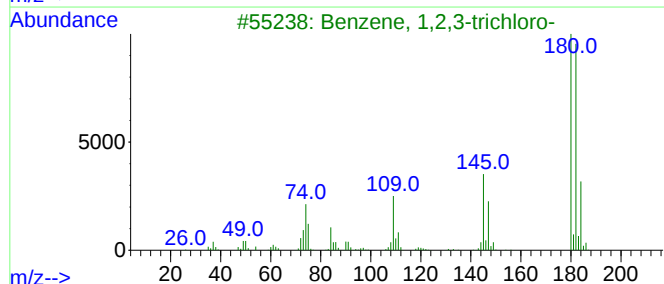
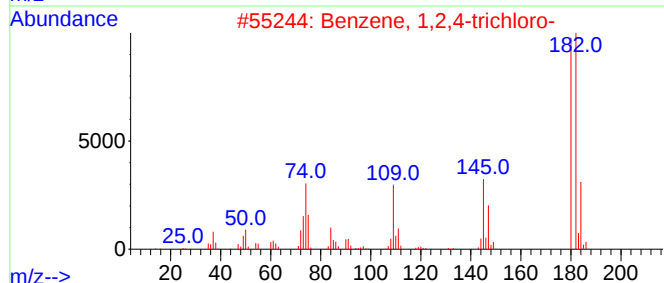
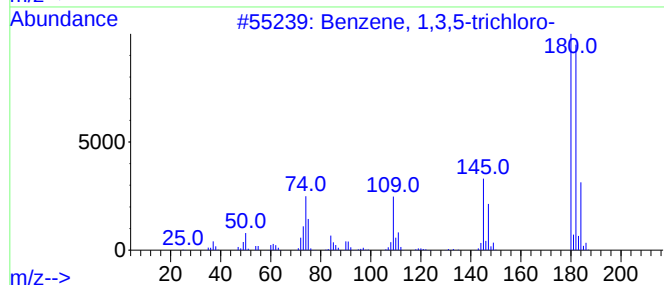
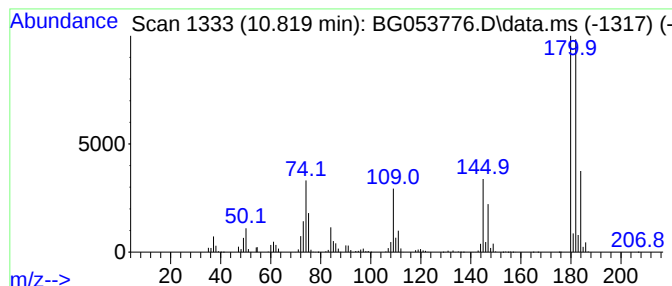
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Benzene, 1,3,5-trichloro- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.819	1990.13 ng/ul	16393800	Naphthalene-d8	10.931

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,3,5-trichloro-	180	C6H3Cl3	000108-70-3	98
2			Benzene, 1,2,4-trichloro-	180	C6H3Cl3	000120-82-1	98
3			Benzene, 1,2,3-trichloro-	180	C6H3Cl3	000087-61-6	98
4			Benzene, 1,3,5-trichloro-	180	C6H3Cl3	000108-70-3	96
5			Benzene, 1,2,3-trichloro-	180	C6H3Cl3	000087-61-6	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG053122\
Data File : BG053776.D
Acq On : 1 Jun 2022 2:35
Operator : CG/JU
Sample : N3039-11
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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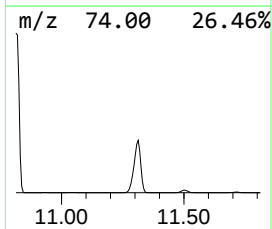
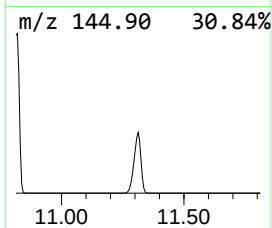
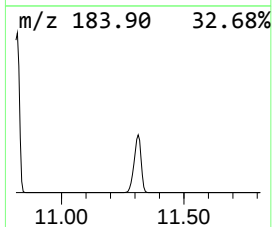
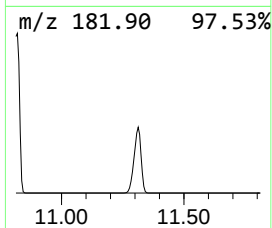
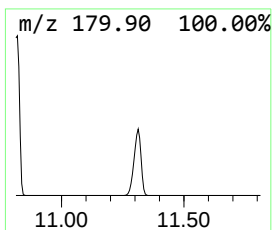
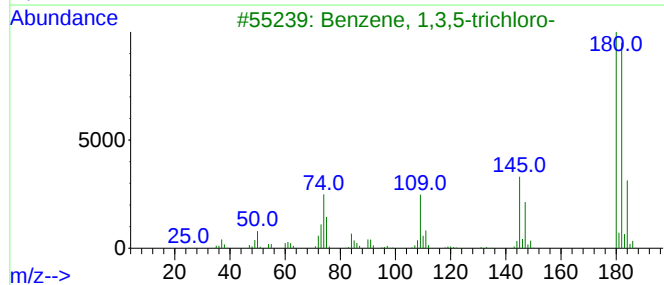
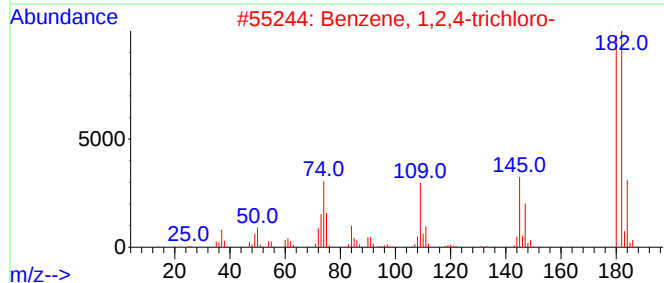
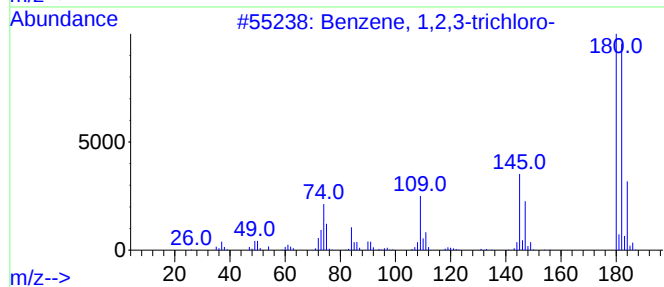
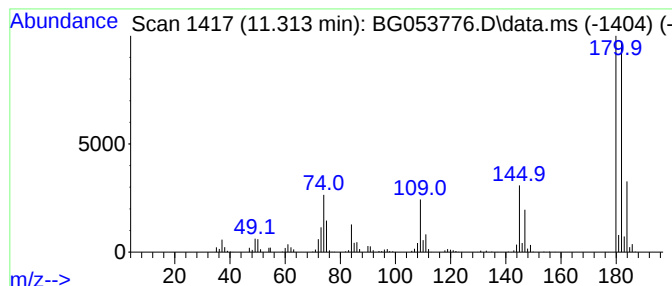
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Benzene, 1,2,3-trichloro- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.313	538.87 ng/ul	4438960	Naphthalene-d8	10.931

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3-trichloro-	180	C6H3Cl3	000087-61-6	99
2			Benzene, 1,2,4-trichloro-	180	C6H3Cl3	000120-82-1	99
3			Benzene, 1,3,5-trichloro-	180	C6H3Cl3	000108-70-3	98
4			Benzene, 1,2,3-trichloro-	180	C6H3Cl3	000087-61-6	97
5			Benzene, 1,3,5-trichloro-	180	C6H3Cl3	000108-70-3	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG053122\
Data File : BG053776.D
Acq On : 1 Jun 2022 2:35
Operator : CG/JU
Sample : N3039-11
Misc :
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Instrument :
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ClientSampleId :
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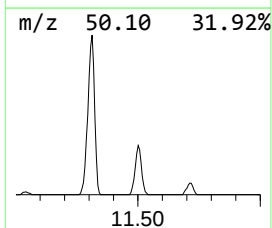
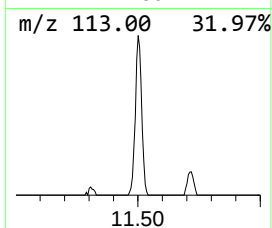
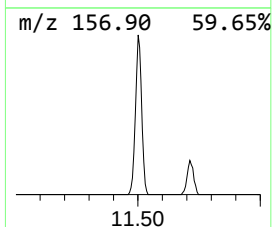
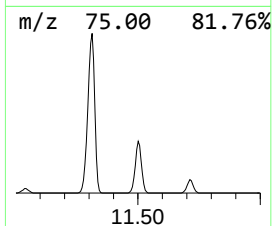
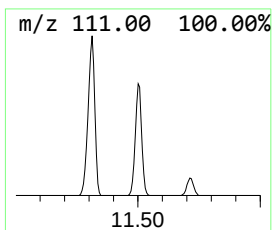
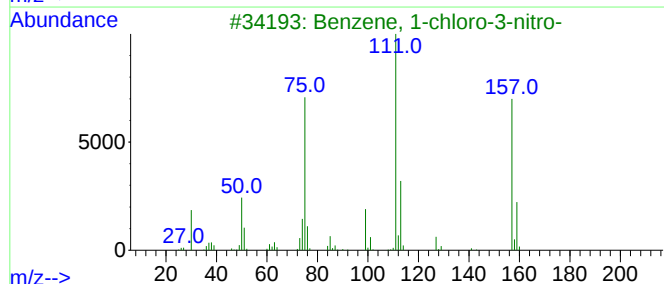
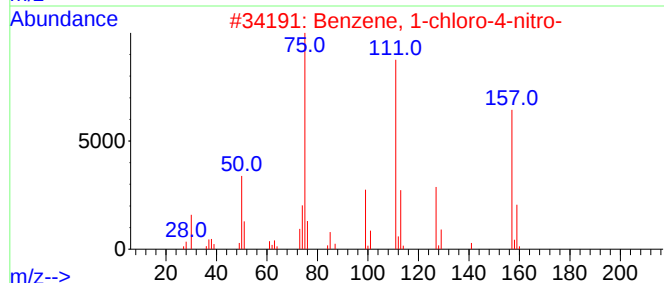
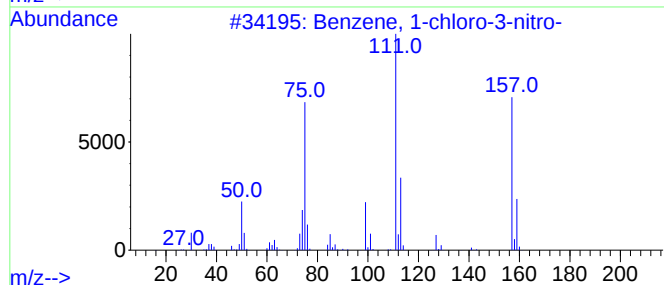
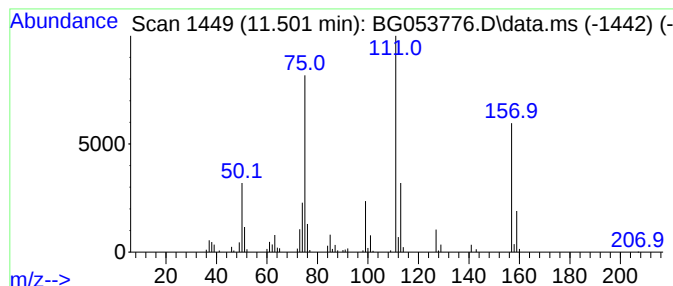
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Benzene, 1-chloro-3-nitro- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.501	29.25 ng/ul	240926	Naphthalene-d8	10.931

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-chloro-3-nitro-	157	C6H4ClNO2	000121-73-3	96
2			Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	94
3			Benzene, 1-chloro-3-nitro-	157	C6H4ClNO2	000121-73-3	93
4			Benzene, 1-chloro-3-nitro-	157	C6H4ClNO2	000121-73-3	93
5			Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG053122\
Data File : BG053776.D
Acq On : 1 Jun 2022 2:35
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Sample : N3039-11
Misc :
ALS Vial : 16 Sample Multiplier: 1

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ClientSampleId :
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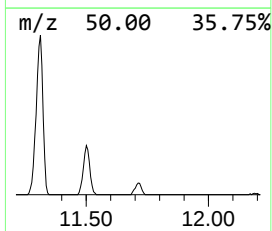
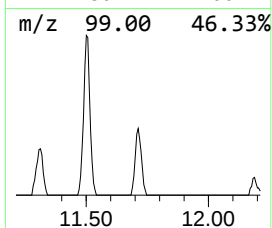
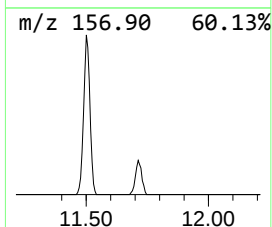
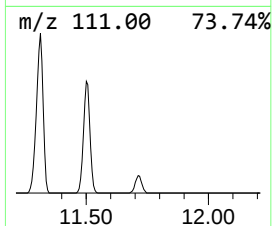
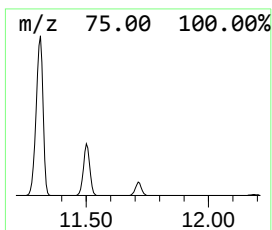
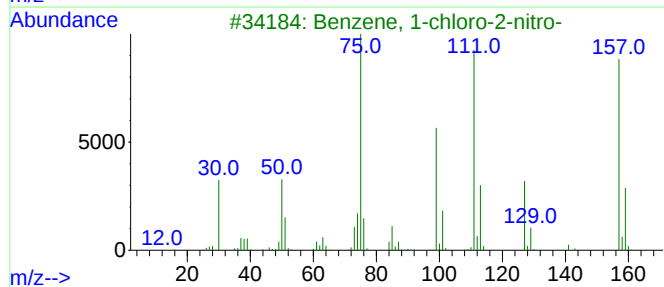
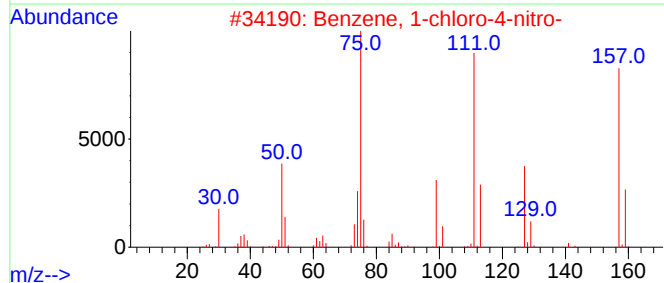
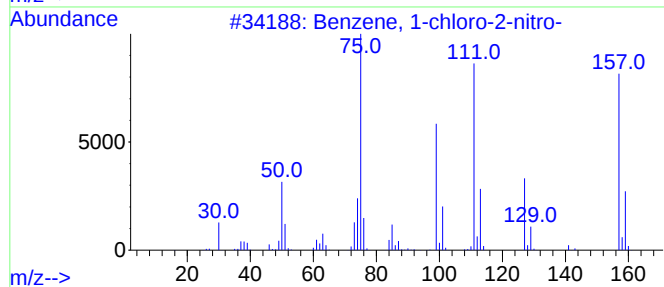
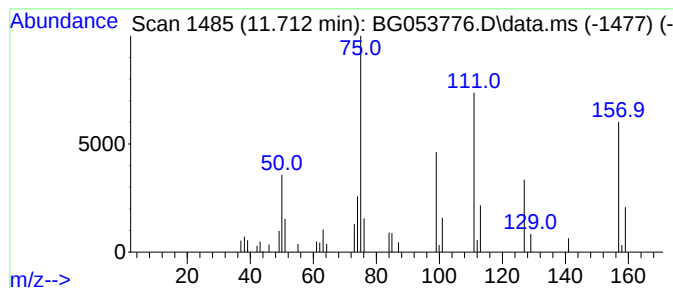
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 Benzene, 1-chloro-2-nitro- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.712	7.07 ng/ul	58277	Naphthalene-d8	10.931

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-chloro-2-nitro-	157	C6H4ClNO2	000088-73-3	98
2			Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	96
3			Benzene, 1-chloro-2-nitro-	157	C6H4ClNO2	000088-73-3	95
4			Benzene, 1-chloro-2-nitro-	157	C6H4ClNO2	000088-73-3	93
5			Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG053122\
Data File : BG053776.D
Acq On : 1 Jun 2022 2:35
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Sample : N3039-11
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ALS Vial : 16 Sample Multiplier: 1

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ClientSampleId :
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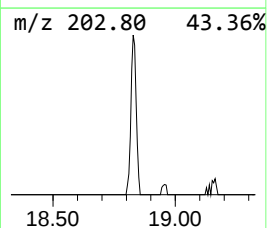
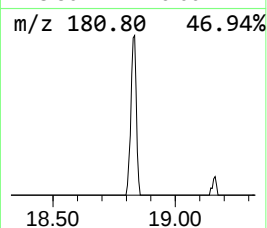
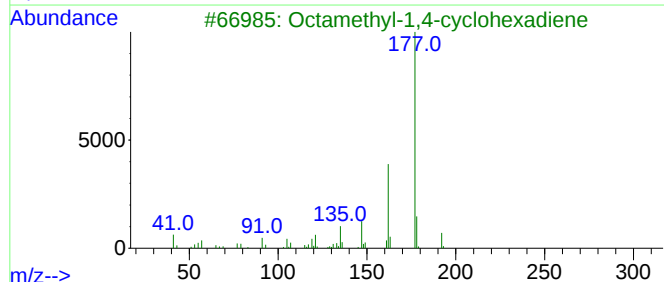
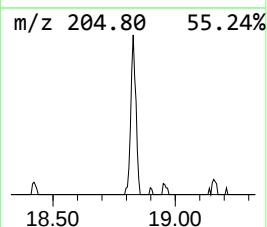
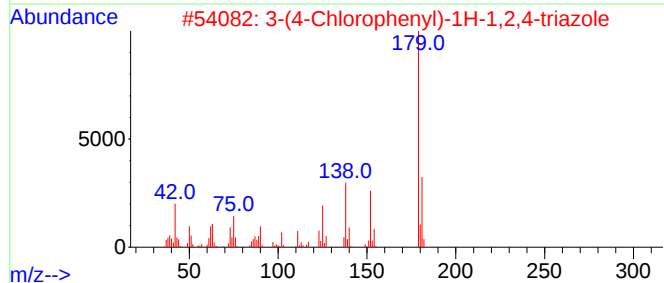
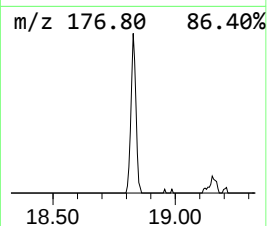
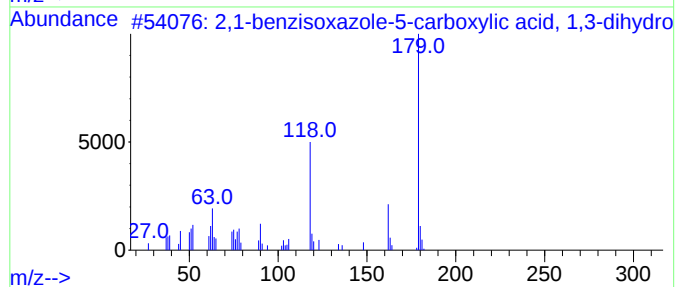
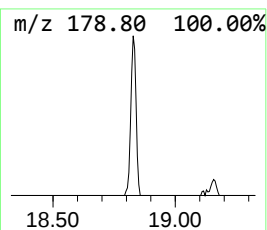
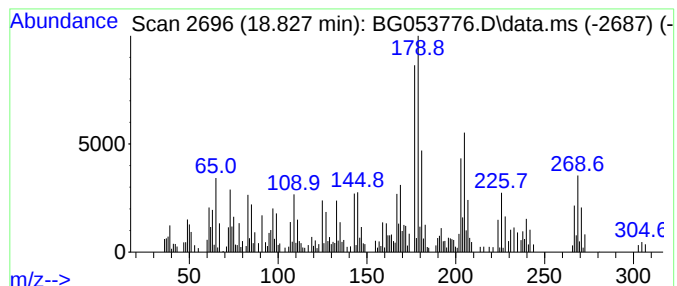
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 unknown-01 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.828	13.68 ng/ul	208265	Phenanthrene-d10	17.464

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,1-benzisoxazole-5-carboxylic a...	179	C8H5N04	1000404-80-3	35		
2	3-(4-Chlorophenyl)-1H-1,2,4-tria...	179	C8H6ClN3	023195-59-7	30		
3	Octamethyl-1,4-cyclohexadiene	192	C14H24	172684-93-4	25		
4	6-Chloro-1,2,3,4-tetrahydrocarba...	205	C12H12ClN	036684-65-8	25		
5	2,4,6-Trifluoronitrobenzene	177	C6H2F3NO2	000315-14-0	18		



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG053122\
Data File : BG053776.D
Acq On : 1 Jun 2022 2:35
Operator : CG/JU
Sample : N3039-11
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0JR4

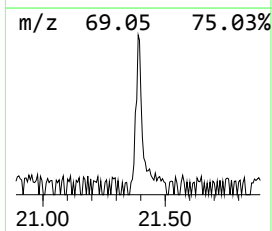
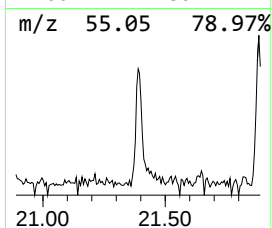
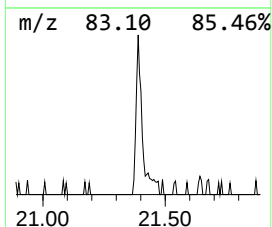
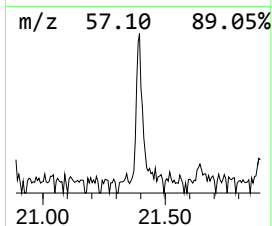
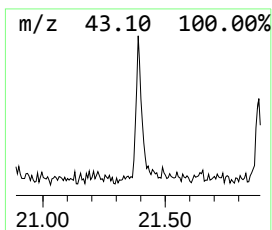
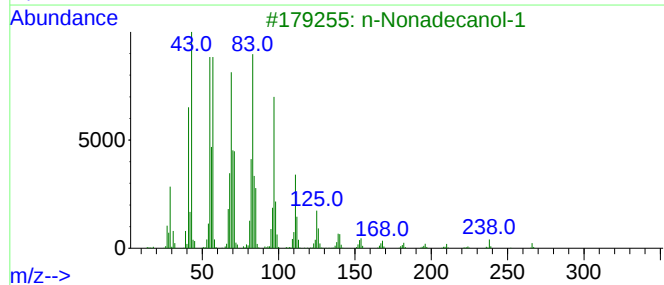
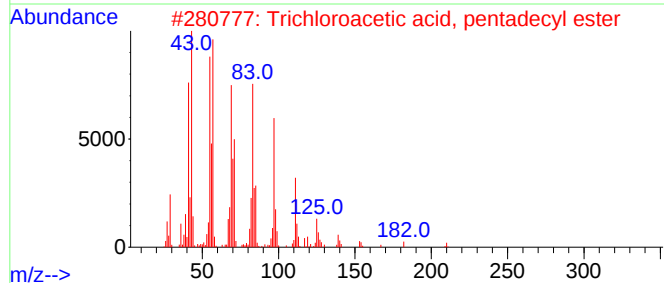
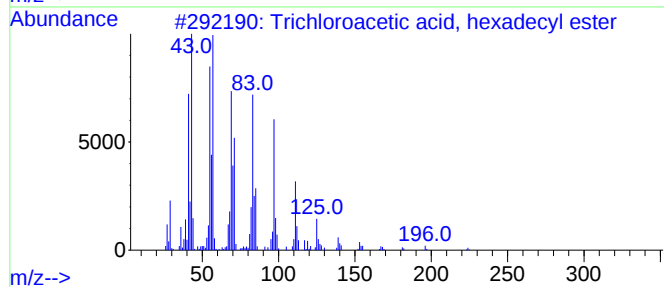
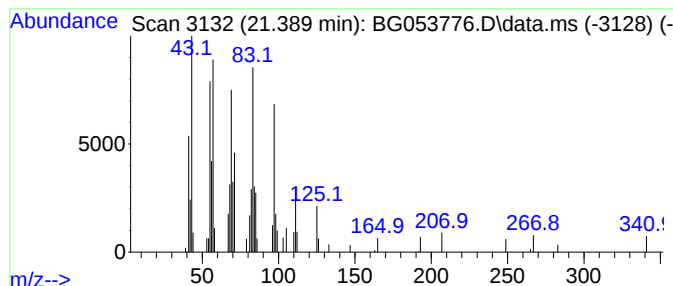
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 Trichloroacetic acid, hexad... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.389	2.16 ng/ul	44311	Chrysene-d12	21.742

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	91
2			Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	91
3			n-Nonadecanol-1	284	C19H40O	001454-84-8	90
4			Trichloroacetic acid, tetradecyl...	358	C16H29Cl3O2	074339-52-9	87
5			1-Hexadecanol	242	C16H34O	036653-82-4	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG053122\
Data File : BG053776.D
Acq On : 1 Jun 2022 2:35
Operator : CG/JU
Sample : N3039-11
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0JR4

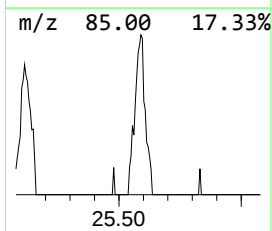
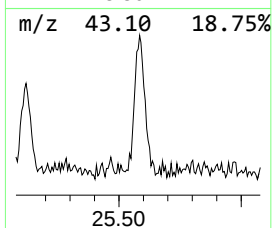
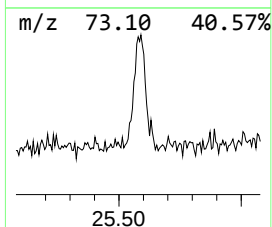
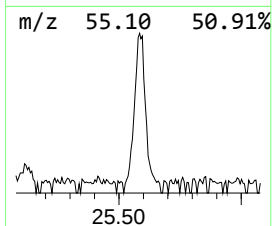
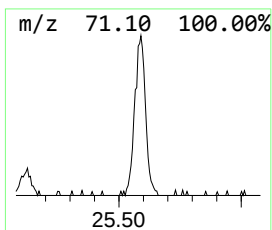
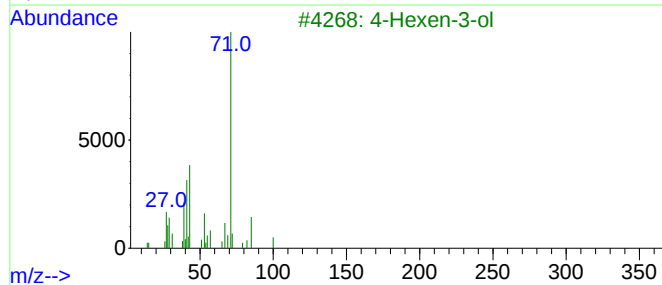
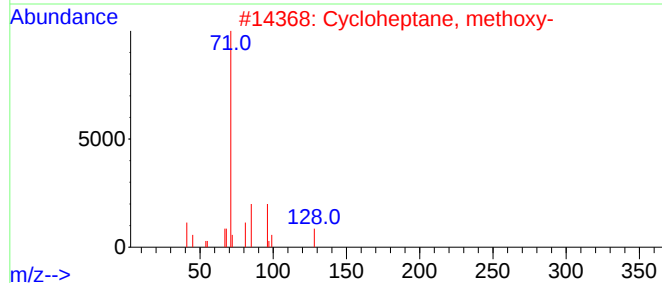
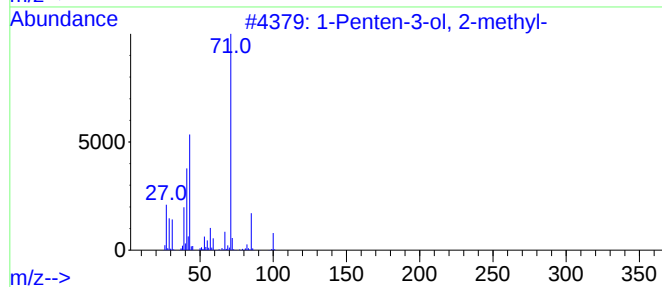
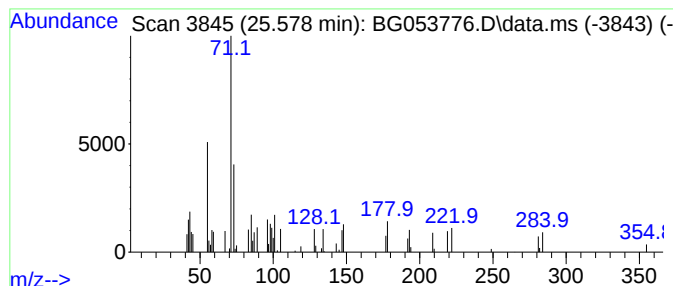
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 unknown-02 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.578	2.71 ng/ul	53495	Perylene-d12	25.014

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Penten-3-ol, 2-methyl-	100	C6H12O	002088-07-5	43
2		Cycloheptane, methoxy-	128	C8H16O	042604-04-6	43
3		4-Hexen-3-ol	100	C6H12O	004798-58-7	43
4		2,2,4-Trimethyl-1,3-pentanediol ...	286	C16H30O4	006846-50-0	32
5		2-Fluorophenethyl alcohol, 2-met...	210	C13H19FO	1000394-96-4	32



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG053122\
 Data File : BG053776.D
 Acq On : 1 Jun 2022 2:35
 Operator : CG/JU
 Sample : N3039-11
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0JR4

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG053122.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
Benzene	3.169	2.5	ng/ul	4191260	1	8.140	33152100	20.0
Benzene, chloro-	5.473	22.9	ng/ul	37938900	1	8.140	33152100	20.0
Benzene, 1,4-di...	8.017	3.9	ng/ul	6522230	1	8.140	33152100	20.0
Benzene, 1,3-di...	8.210	20.0	ng/ul	33152100	1	8.140	33152100	20.0
Benzene, 1,2-di...	8.557	23.9	ng/ul	39591500	1	8.140	33152100	20.0
Benzene, 1-brom...	9.603	17.0	ng/ul	140310	2	10.931	164751	20.0
Benzene, 1-brom...	9.920	10.4	ng/ul	85762	2	10.931	164751	20.0
Benzene, 1,3,5-...	10.819	1990.1	ng/ul	16393800	2	10.931	164751	20.0
Benzene, 1,2,3-...	11.313	538.9	ng/ul	4438960	2	10.931	164751	20.0
Benzene, 1-chlo...	11.501	29.3	ng/ul	240926	2	10.931	164751	20.0
Benzene, 1-chlo...	11.712	7.1	ng/ul	58277	2	10.931	164751	20.0
unknown-01	18.828	13.7	ng/ul	208265	4	17.464	304483	20.0
Trichloroacetic...	21.389	2.2	ng/ul	44311	5	21.742	410800	20.0
unknown-02	25.578	2.7	ng/ul	53495	6	25.014	394558	20.0