

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
 Data File : BP010038.D
 Acq On : 22 Apr 2022 05:06
 Operator : CG/JU
 Sample : N2490-06
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0J23

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_P\Methods\SFAM-EPA-BP042122.M
 Title : SVOA CALIBRATION

Signal : TIC: BP010038.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.411	360	395	399	rBV4	53542107	491420877	100.00%	35.210%
2	6.881	627	645	651	rBV	605410	1163849	0.24%	0.083%
3	7.134	678	688	703	rVB2	283954	775567	0.16%	0.056%
4	7.299	703	716	731	rBV	3250731	9728920	1.98%	0.697%
5	7.516	740	753	766	rVB3	568127	2122254	0.43%	0.152%
6	7.858	799	811	826	rBV	9711013	42791374	8.71%	3.066%
7	8.105	826	853	857	rVB3	49797786	340557275	69.30%	24.400%
8	8.452	880	912	916	rBV3	50278294	348678584	70.95%	24.982%
9	8.658	937	947	959	rVV	484214	789473	0.16%	0.057%
10	9.111	1017	1024	1029	rBV	771072	1274470	0.26%	0.091%
11	9.452	1073	1082	1090	rBV	322270	554731	0.11%	0.040%
12	9.781	1129	1138	1143	rBV	4213023	7234679	1.47%	0.518%
13	9.834	1143	1147	1156	rVB	675922	1055726	0.21%	0.076%
14	10.369	1232	1238	1245	rBV	854233	1546376	0.31%	0.111%
15	10.440	1245	1250	1257	rVB	661812	1136515	0.23%	0.081%
16	10.534	1257	1266	1273	rBV	1338151	2420058	0.49%	0.173%
17	10.634	1273	1283	1297	rVB	23459590	46262468	9.41%	3.315%
18	10.757	1297	1304	1313	rVB	796062	1255266	0.26%	0.090%
19	10.893	1313	1327	1331	rBV	17854002	43265710	8.80%	3.100%
20	10.928	1331	1333	1343	rVB	2060865	2337056	0.48%	0.167%
21	11.146	1362	1370	1377	rBV	6410163	10892270	2.22%	0.780%
22	12.775	1639	1647	1656	rVB	922494	1561817	0.32%	0.112%
23	13.004	1679	1686	1693	rVV2	435870	783258	0.16%	0.056%
24	13.581	1775	1784	1791	rBV	900516	1403417	0.29%	0.101%
25	13.975	1845	1851	1865	rVB	1794102	2659739	0.54%	0.191%
26	14.263	1894	1900	1909	rVB	2060487	2935544	0.60%	0.210%
27	14.569	1945	1952	1964	rBV	1206738	1830155	0.37%	0.131%
28	15.563	2115	2121	2132	rVB	2626741	3803380	0.77%	0.273%
29	15.675	2132	2140	2150	rBV	846520	1235556	0.25%	0.089%
30	17.322	2413	2420	2430	rBV	1588740	2292906	0.47%	0.164%
31	17.422	2430	2437	2447	rBV	3139807	4346319	0.88%	0.311%
32	18.663	2639	2648	2656	rBV	493672	683865	0.14%	0.049%
33	19.651	2810	2816	2830	rBV	4094274	5361638	1.09%	0.384%
34	21.398	3108	3113	3127	rBV	1797537	2619453	0.53%	0.188%
35	23.698	3495	3504	3511	rBV2	2138738	4580413	0.93%	0.328%
36	23.851	3523	3530	3545	rVB2	1123370	2341711	0.48%	0.168%

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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_P\Methods\SFAM-EPA-BP042122.M
Title : SVOA CALIBRATION

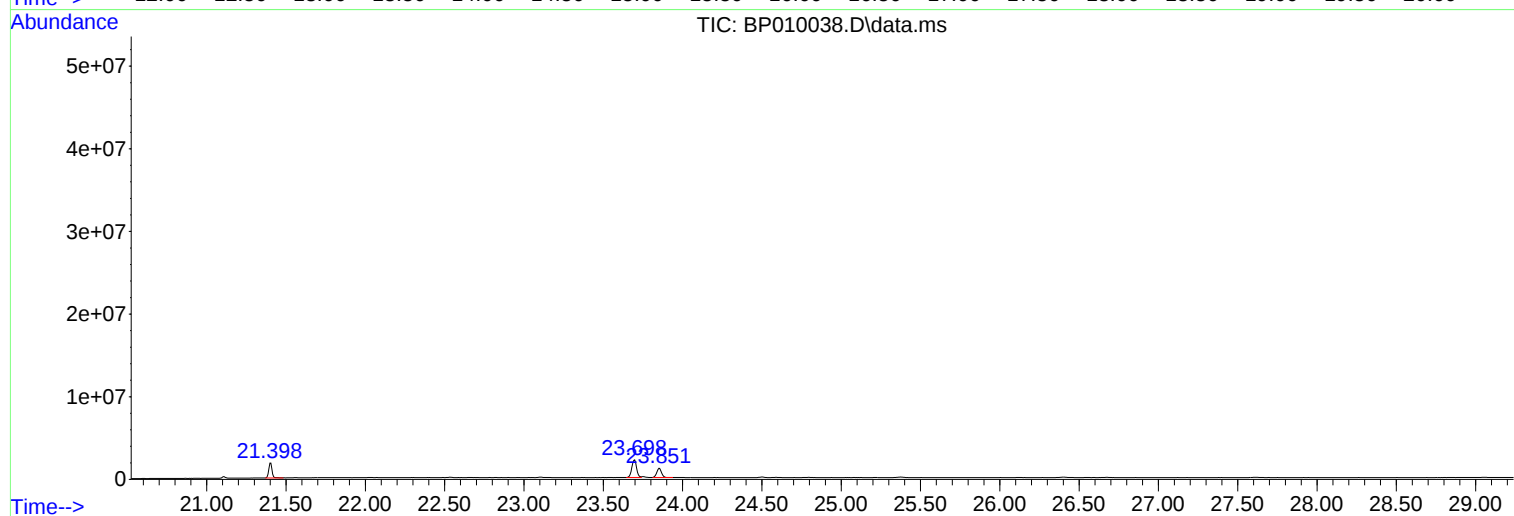
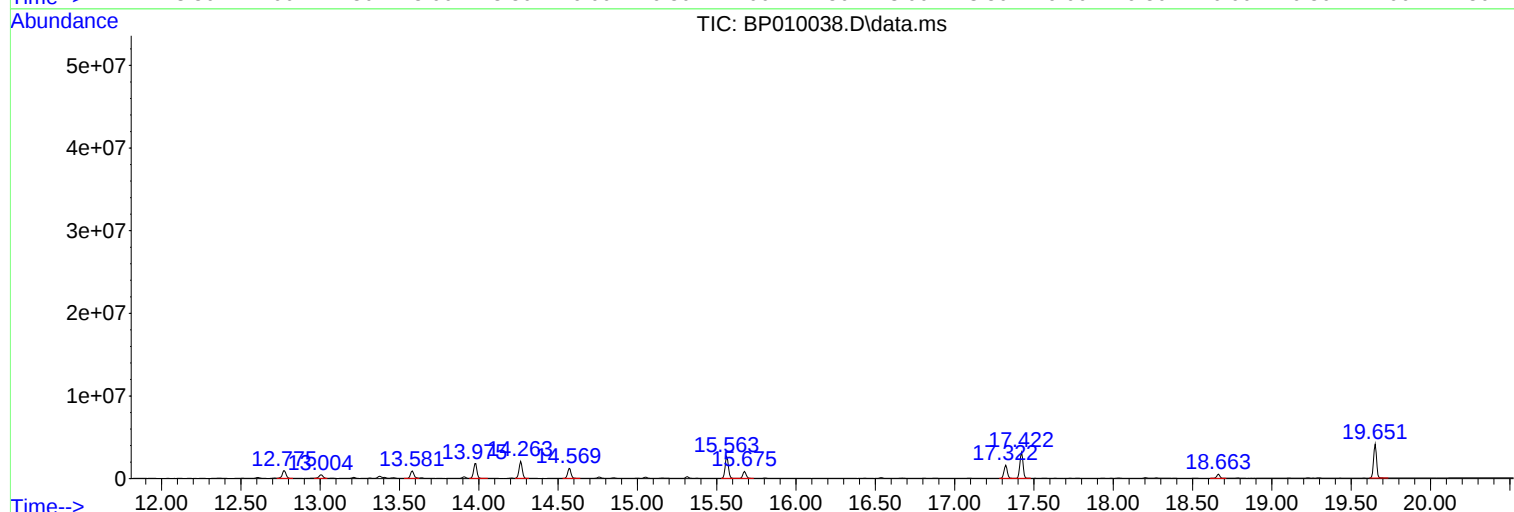
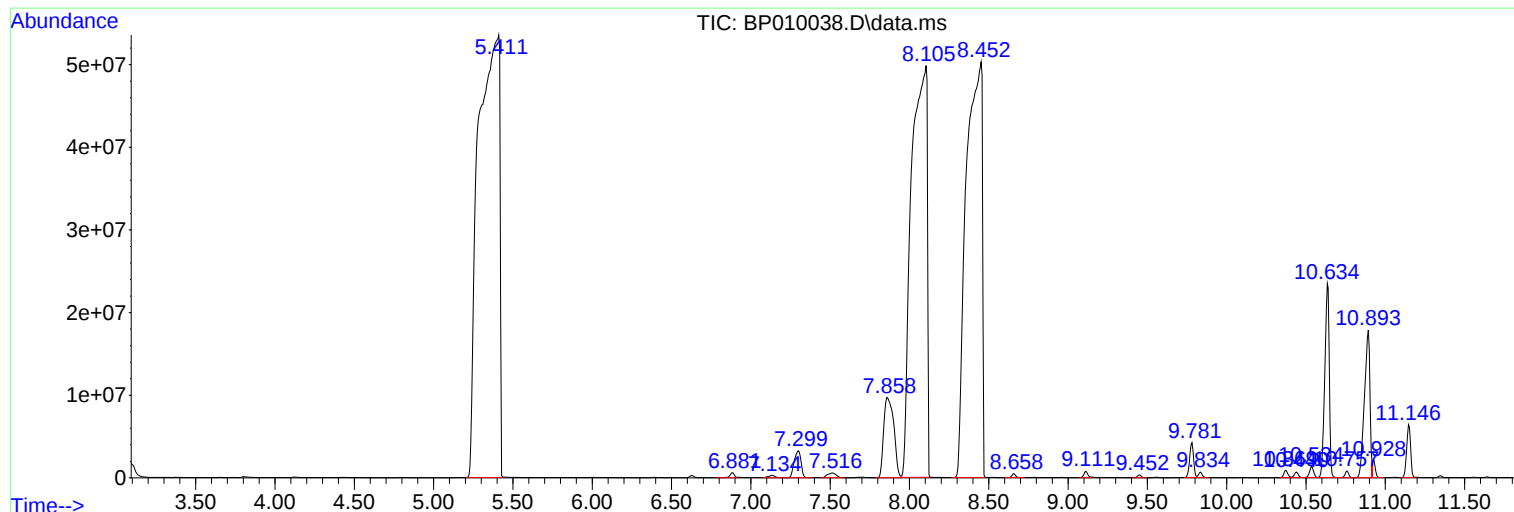
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

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



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Instrument :
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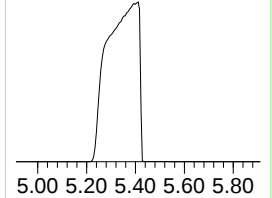
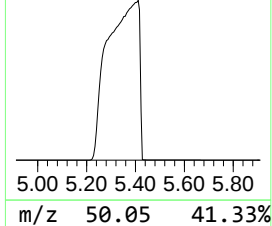
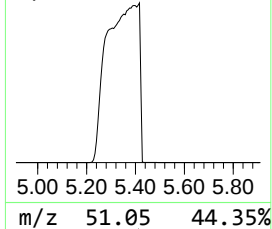
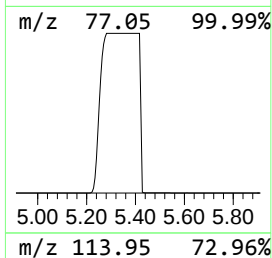
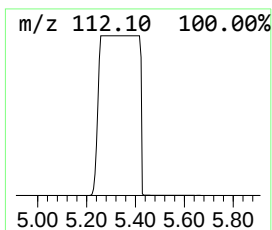
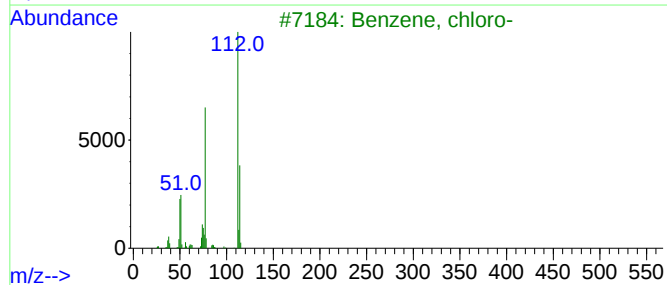
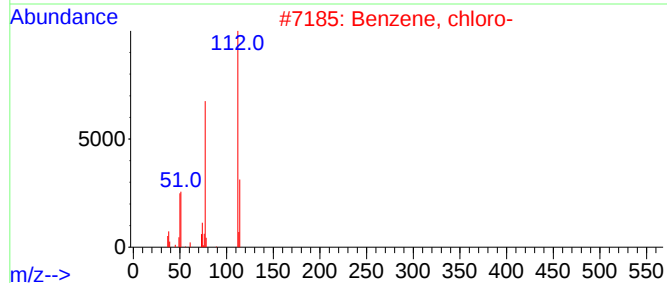
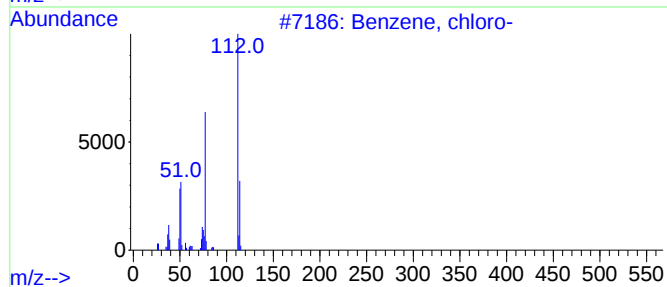
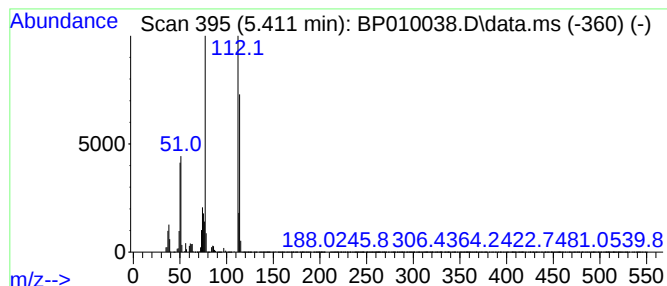
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Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzene, chloro- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.411	28.86 ng/ul	491421000	1,4-Dichlorobenzene-d4	8.022

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, chloro-	112	C6H5Cl	000108-90-7	95
2			Benzene, chloro-	112	C6H5Cl	000108-90-7	94
3			Benzene, chloro-	112	C6H5Cl	000108-90-7	94
4			Benzene, chloro-	112	C6H5Cl	000108-90-7	50
5			Benzene, chloro-	112	C6H5Cl	000108-90-7	38



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Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0J23

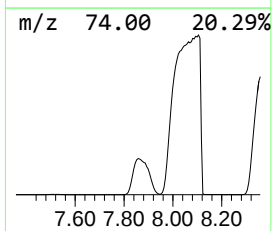
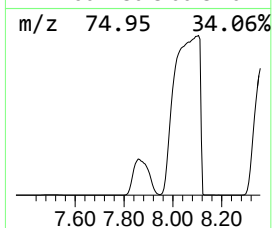
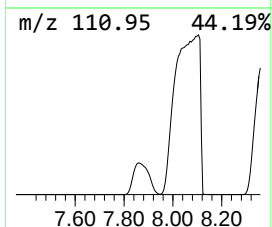
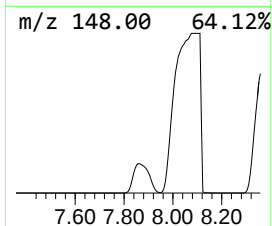
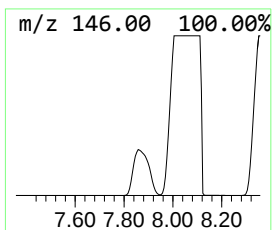
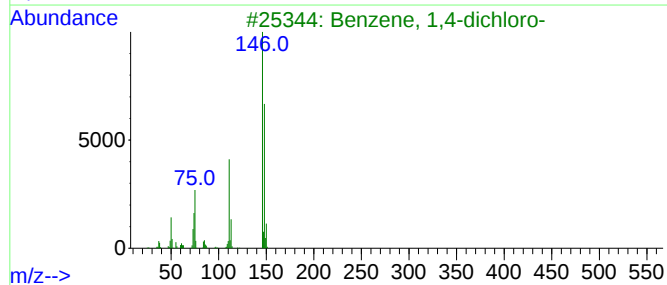
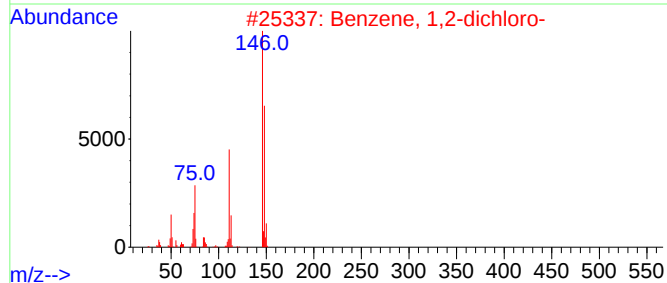
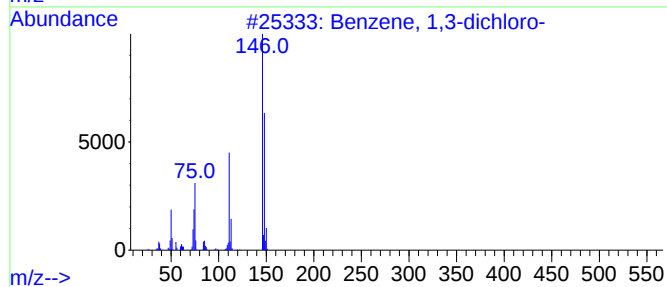
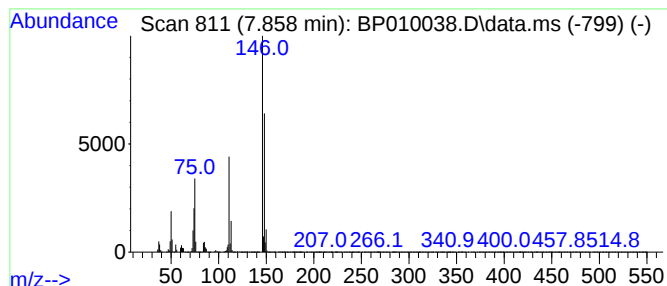
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzene, 1,3-dichloro- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.858	2.51 ng/ul	42791400	1,4-Dichlorobenzene-d4	8.022

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	98
2			Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	98
3			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	98
4			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	96
5			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	96



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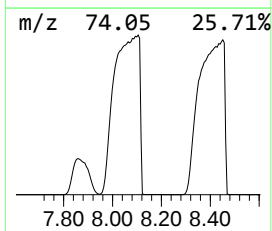
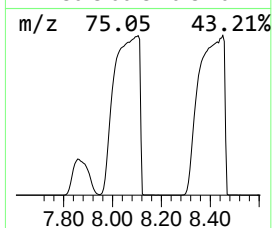
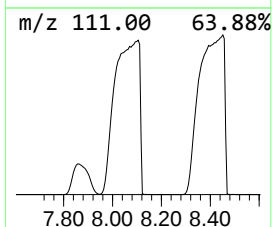
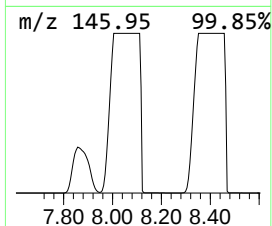
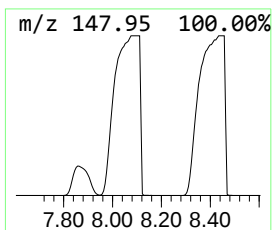
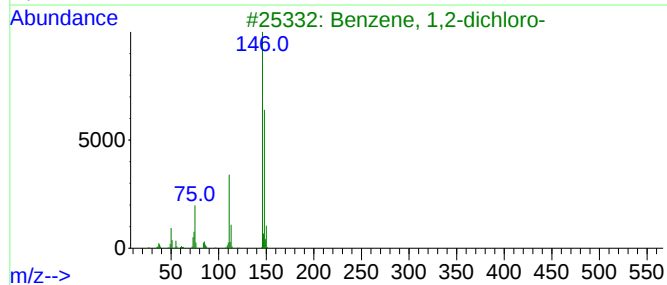
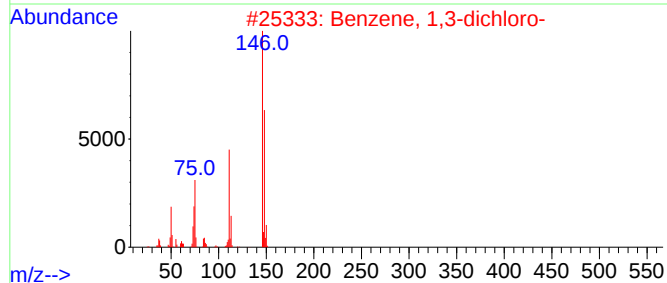
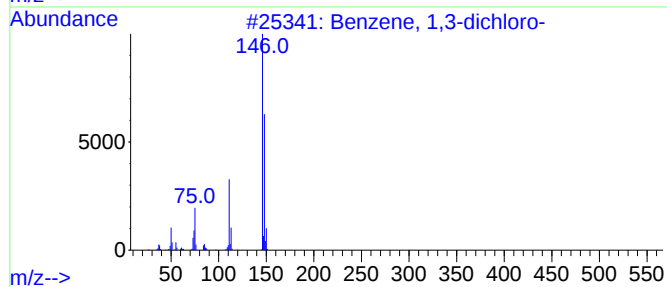
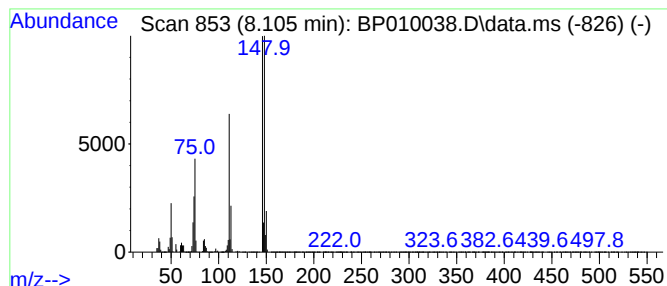
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Benzene, 1,2-dichloro- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.105	20.00 ng/ul	340557000	1,4-Dichlorobenzene-d4	8.022

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	95
2			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	95
3			Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	93
4			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	93
5			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	93



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Instrument :
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ClientSampleId :
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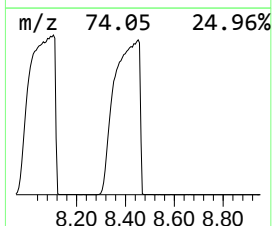
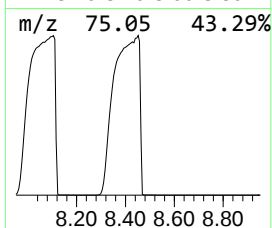
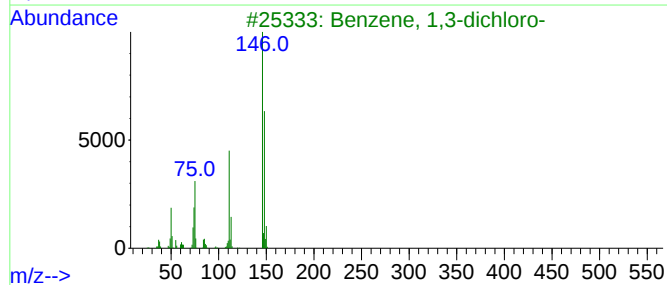
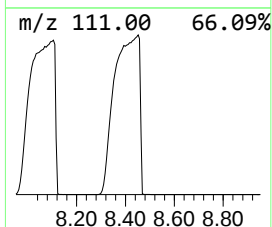
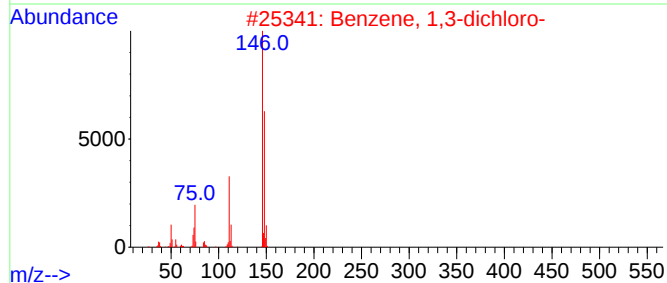
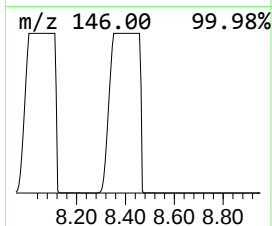
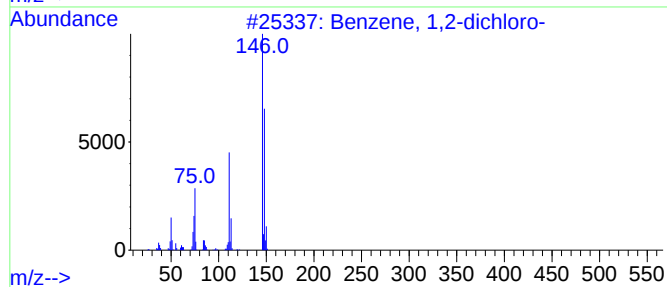
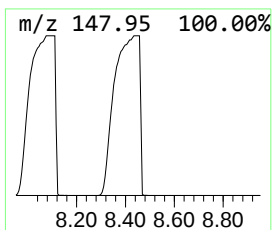
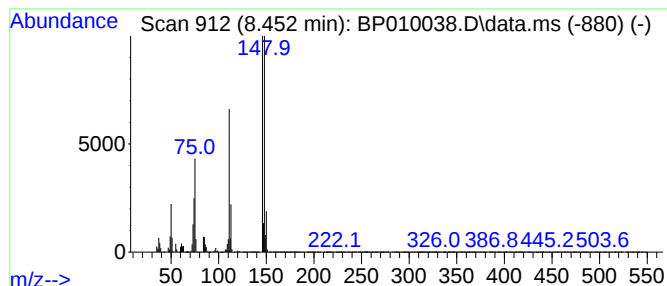
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.452	20.48 ng/ul	348679000	1,4-Dichlorobenzene-d4	8.022

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	94
2			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	93
3			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	93
4			Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	93
5			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	93



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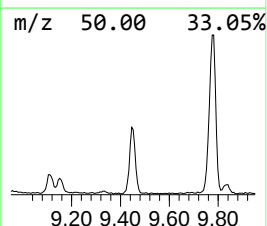
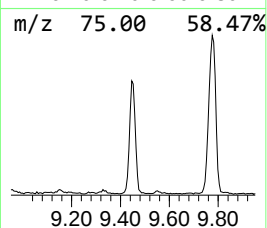
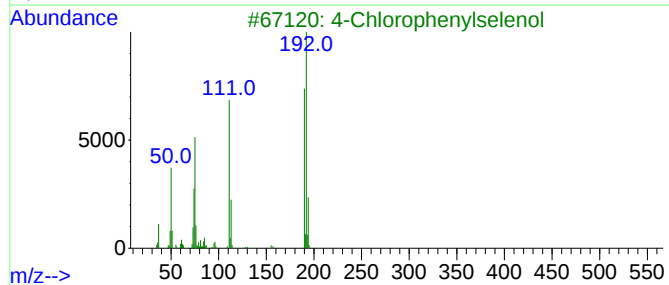
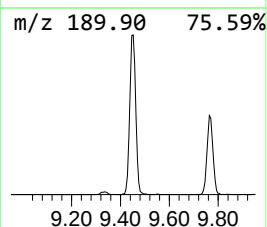
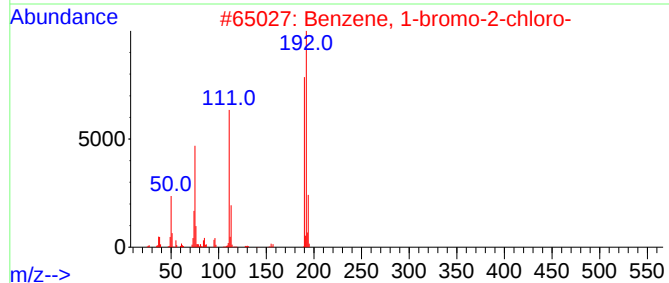
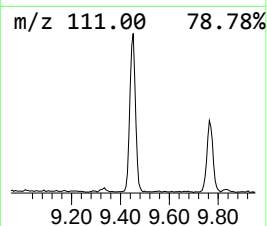
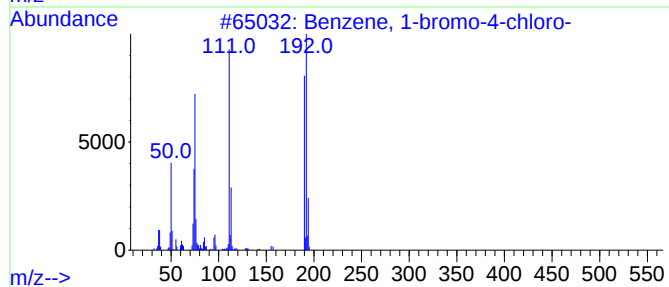
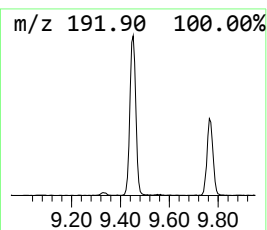
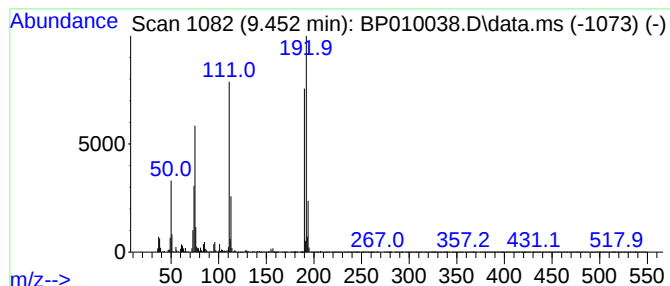
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

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

Peak Number 5 Benzene, 1-bromo-4-chloro- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.452	8.84 ng/ul	554731	Naphthalene-d8	10.758

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	99
2			Benzene, 1-bromo-2-chloro-	190	C6H4BrCl	000694-80-4	98
3			4-Chlorophenylselenol	192	C6H5ClSe	016645-10-6	98
4			Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	98
5			Benzene, 1-bromo-3-chloro-	190	C6H4BrCl	000108-37-2	98



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010038.D
Acq On : 22 Apr 2022 05:06
Operator : CG/JU
Sample : N2490-06
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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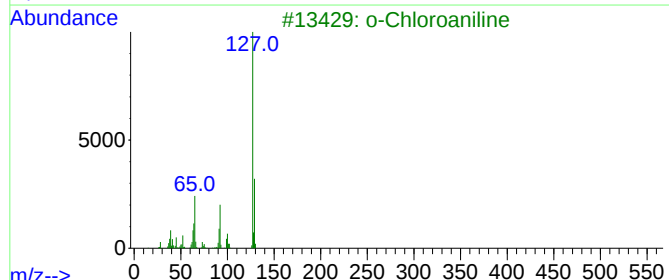
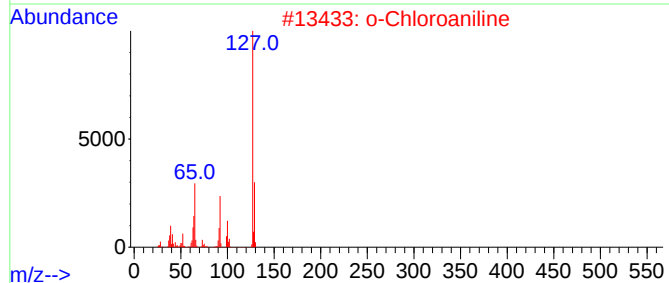
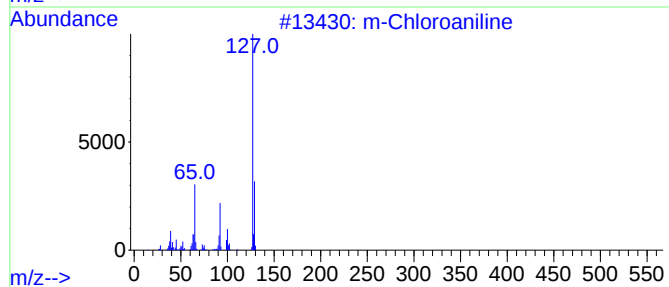
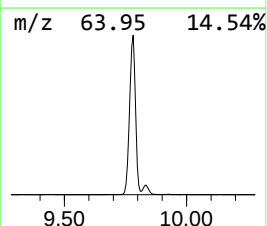
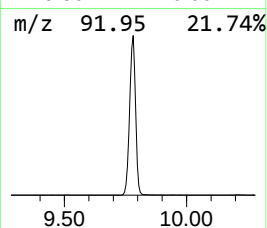
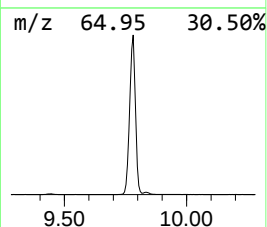
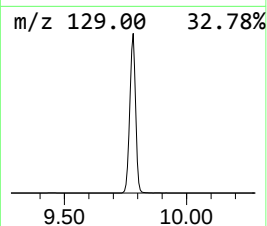
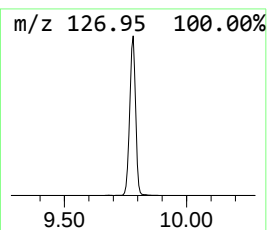
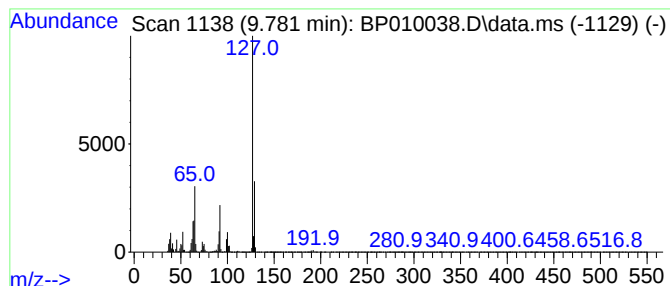
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 m-Chloroaniline Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.781	115.27 ng/ul	7234680	Naphthalene-d8	10.758

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			m-Chloroaniline	127	C6H6ClN	000108-42-9	97
2			o-Chloroaniline	127	C6H6ClN	000095-51-2	96
3			o-Chloroaniline	127	C6H6ClN	000095-51-2	96
4			o-Chloroaniline	127	C6H6ClN	000095-51-2	95
5			m-Chloroaniline	127	C6H6ClN	000108-42-9	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010038.D
Acq On : 22 Apr 2022 05:06
Operator : CG/JU
Sample : N2490-06
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Instrument :
BNA_P
ClientSampleId :
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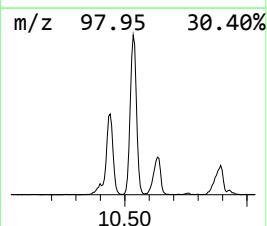
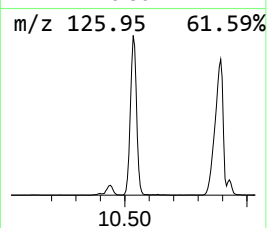
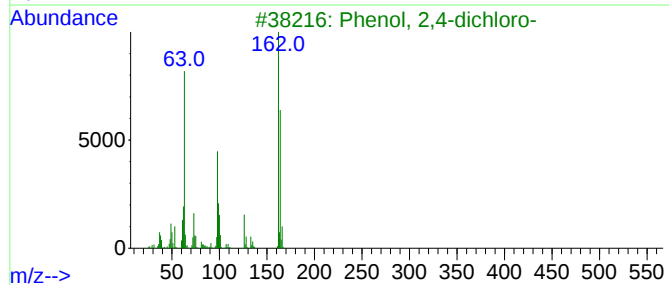
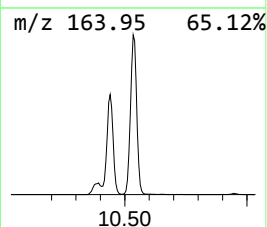
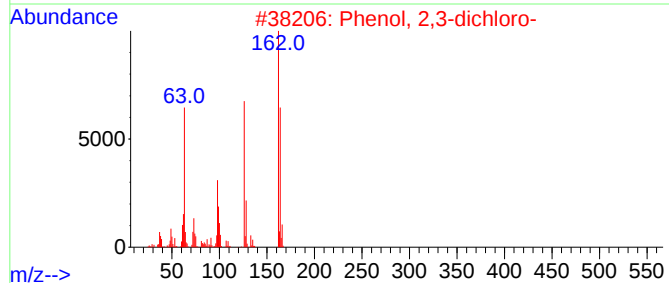
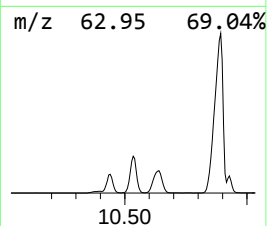
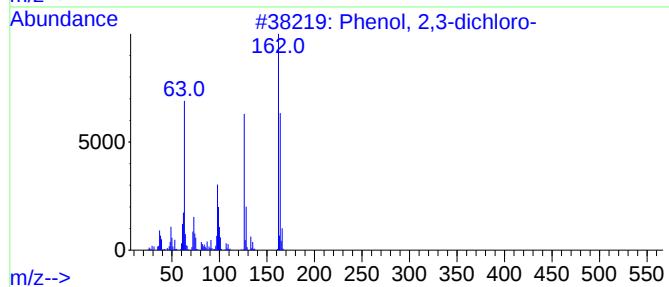
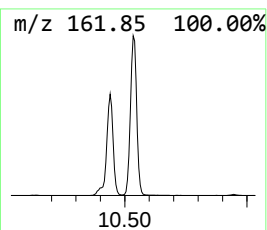
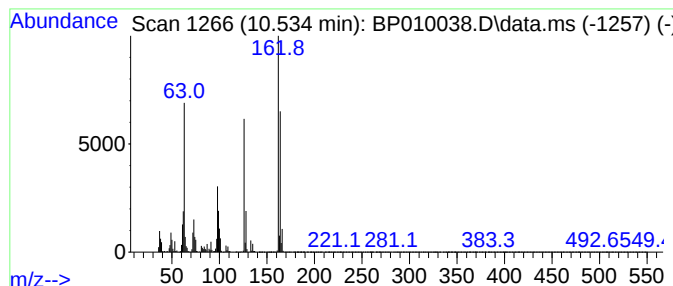
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Phenol, 2,3-dichloro- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.534	38.56 ng/ul	2420060	Naphthalene-d8	10.758

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 2,3-dichloro-	162	C6H4Cl2O	000576-24-9	99
2			Phenol, 2,3-dichloro-	162	C6H4Cl2O	000576-24-9	98
3			Phenol, 2,4-dichloro-	162	C6H4Cl2O	000120-83-2	98
4			Phenol, 2,3-dichloro-	162	C6H4Cl2O	000576-24-9	97
5			Phenol, 2,3-dichloro-	162	C6H4Cl2O	000576-24-9	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010038.D
Acq On : 22 Apr 2022 05:06
Operator : CG/JU
Sample : N2490-06
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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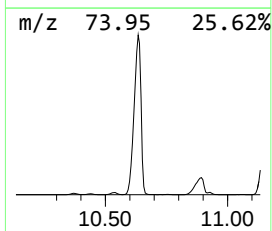
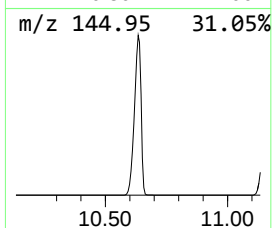
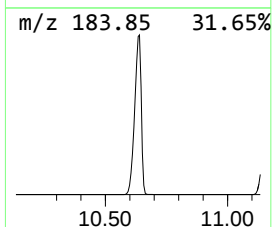
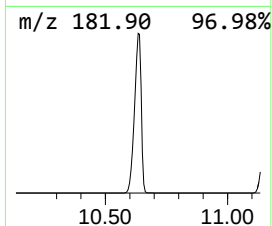
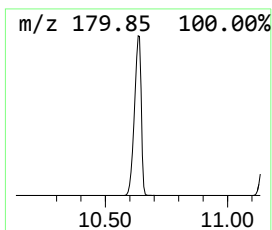
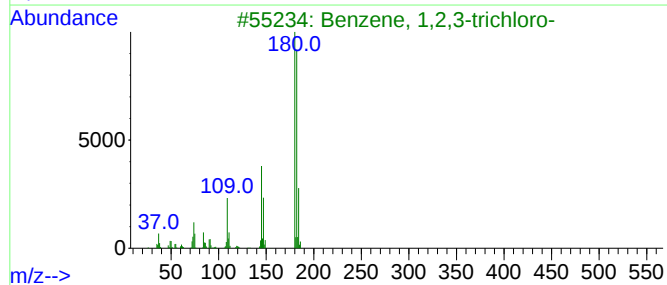
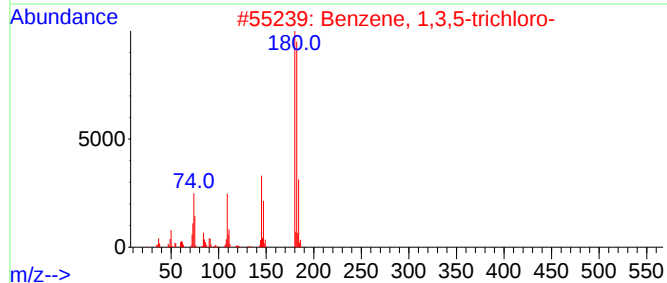
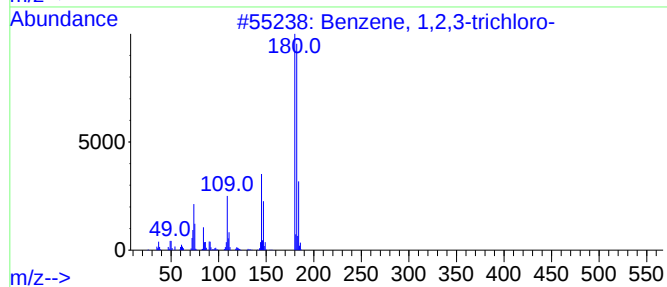
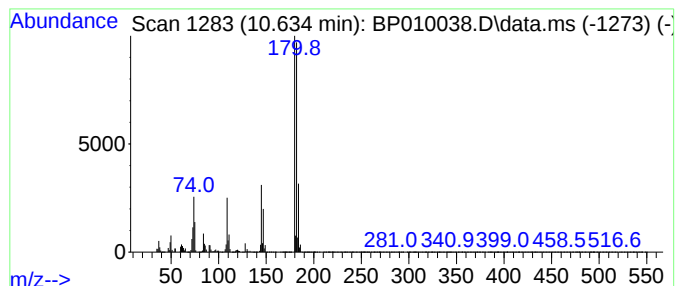
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Benzene, 1,2,3-trichloro- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.634	737.09 ng/ul	46262500	Naphthalene-d8	10.758

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010038.D
Acq On : 22 Apr 2022 05:06
Operator : CG/JU
Sample : N2490-06
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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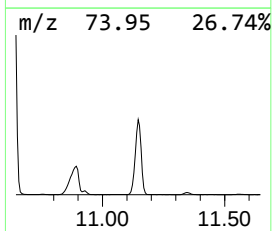
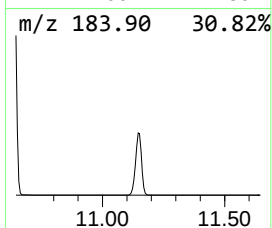
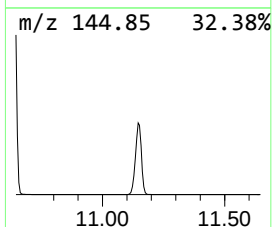
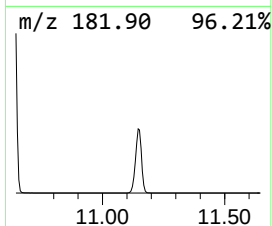
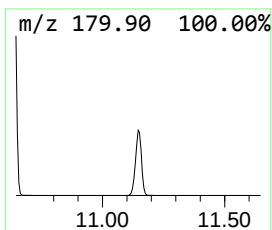
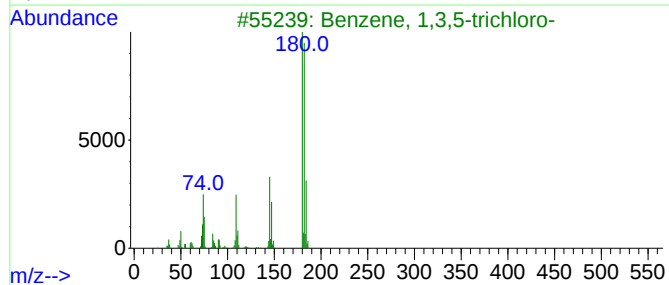
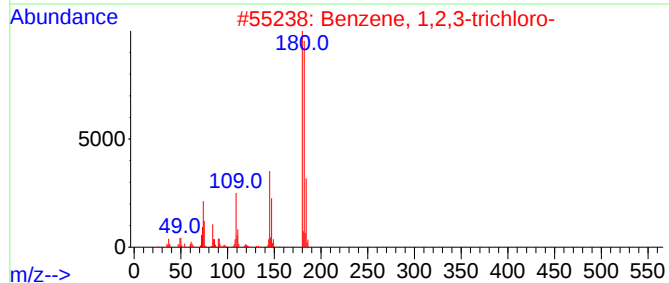
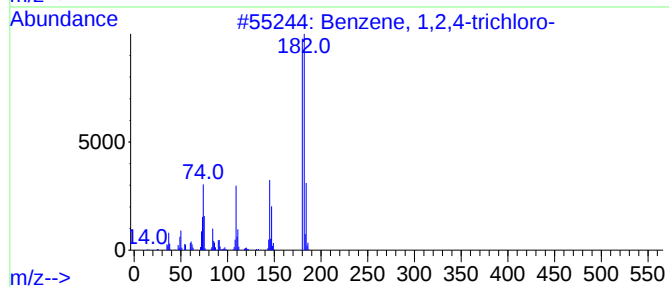
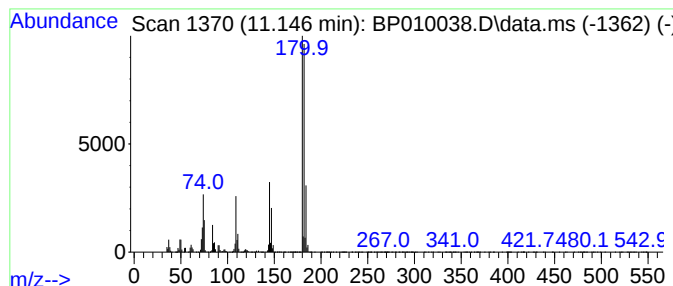
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Benzene, 1,2,4-trichloro- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.146	173.54 ng/ul	10892300	Naphthalene-d8	10.758

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010038.D
Acq On : 22 Apr 2022 05:06
Operator : CG/JU
Sample : N2490-06
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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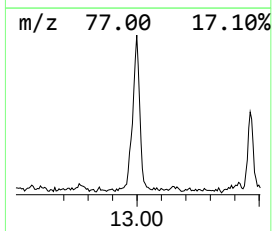
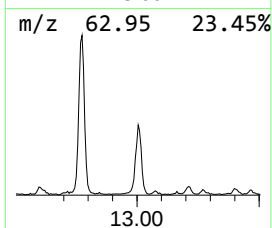
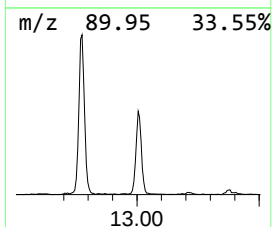
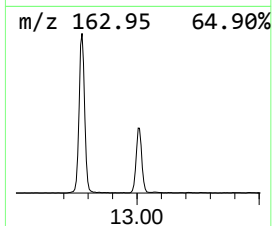
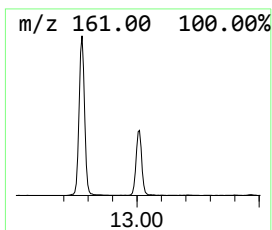
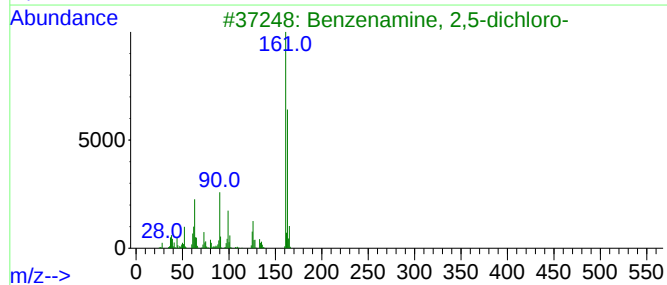
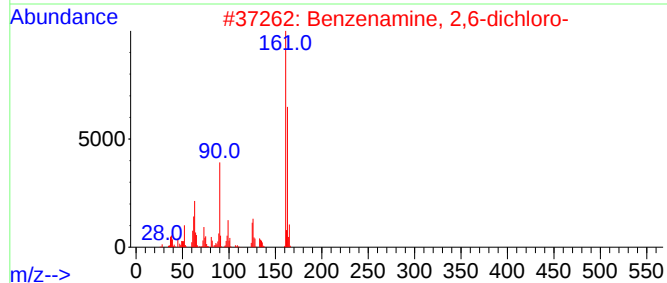
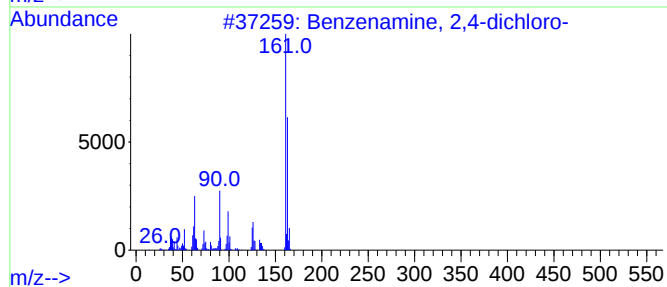
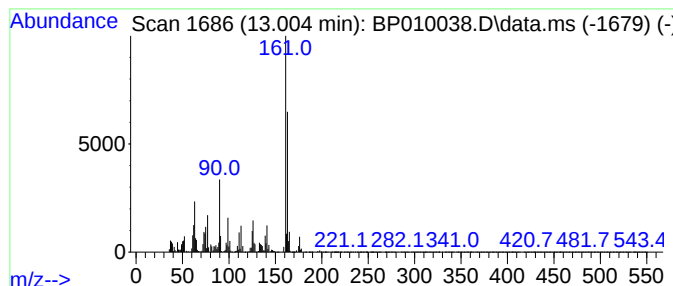
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Benzenamine, 2,4-dichloro- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.004	8.56 ng/ul	783258	Acenaphthene-d10	14.569

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzenamine, 2,4-dichloro-	161	C6H5Cl2N	000554-00-7	98
2			Benzenamine, 2,6-dichloro-	161	C6H5Cl2N	000608-31-1	98
3			Benzenamine, 2,5-dichloro-	161	C6H5Cl2N	000095-82-9	98
4			Benzenamine, 2,3-dichloro-	161	C6H5Cl2N	000608-27-5	98
5			Benzenamine, 2,5-dichloro-	161	C6H5Cl2N	000095-82-9	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010038.D
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Misc :
ALS Vial : 34 Sample Multiplier: 1

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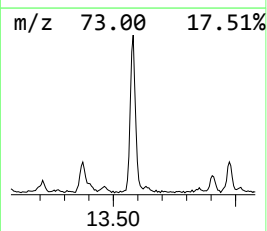
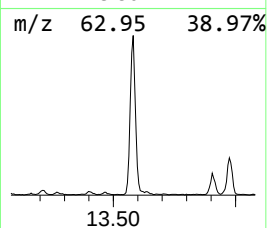
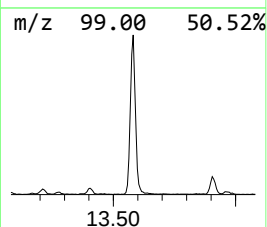
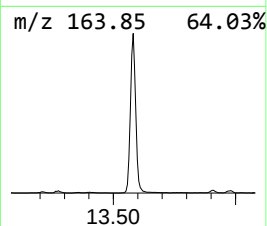
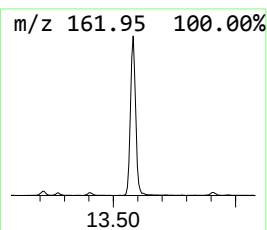
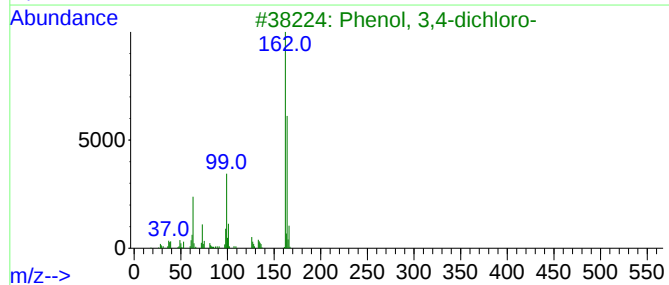
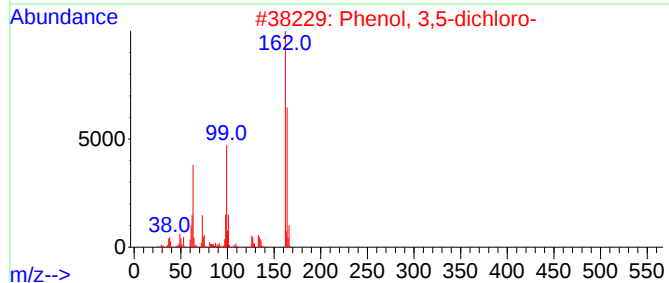
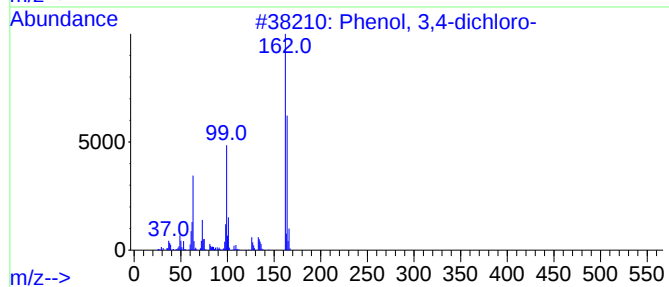
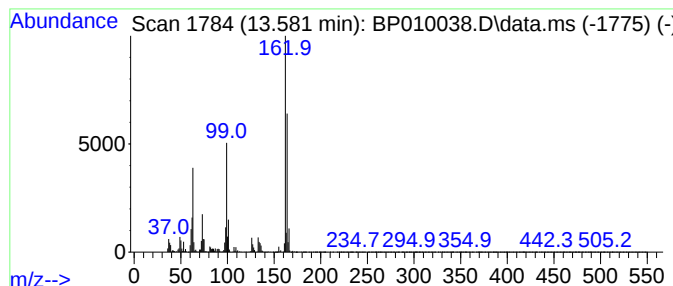
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 Phenol, 3,4-dichloro- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.581	15.34 ng/ul	1403420	Acenaphthene-d10	14.569

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 3,4-dichloro-	162	C6H4Cl2O	000095-77-2	98
2			Phenol, 3,5-dichloro-	162	C6H4Cl2O	000591-35-5	98
3			Phenol, 3,4-dichloro-	162	C6H4Cl2O	000095-77-2	97
4			Phenol, 3,4-dichloro-	162	C6H4Cl2O	000095-77-2	97
5			Phenol, 3,5-dichloro-	162	C6H4Cl2O	000591-35-5	97



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010038.D
Acq On : 22 Apr 2022 05:06
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Sample : N2490-06
Misc :
ALS Vial : 34 Sample Multiplier: 1

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ClientSampleId :
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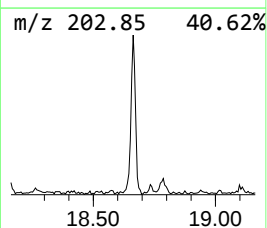
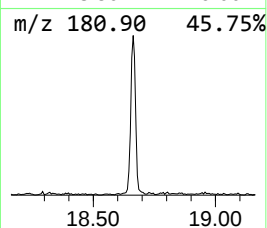
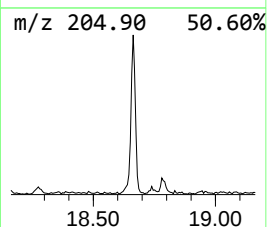
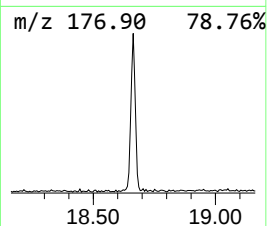
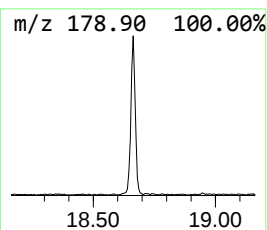
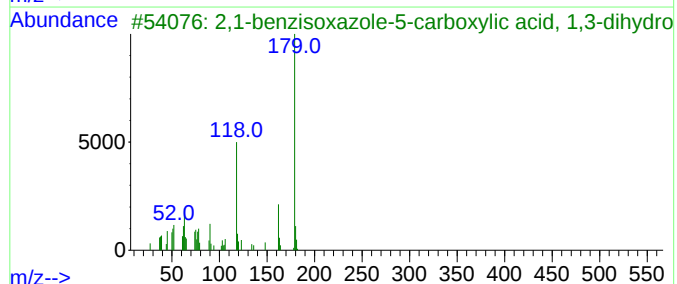
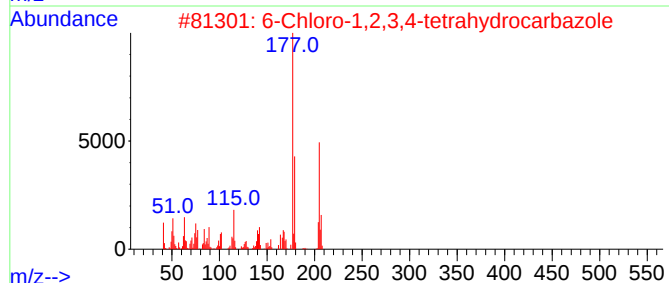
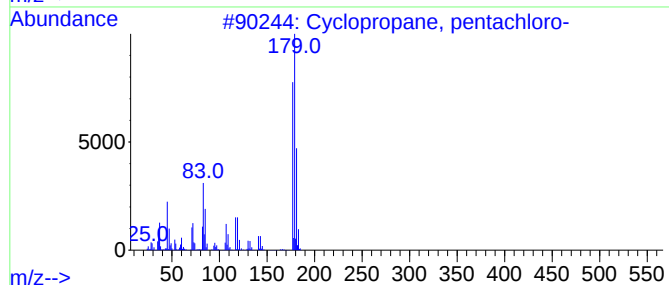
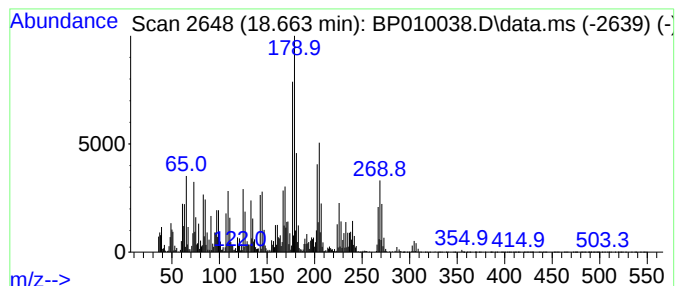
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

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

Peak Number 12 unknown-01 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.663	5.97 ng/ul	683865	Phenanthrene-d10	17.322

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopropane, pentachloro-	212	C3HCl5	006262-51-7	43
2			6-Chloro-1,2,3,4-tetrahydrocarba...	205	C12H12ClN	036684-65-8	38
3			2,1-benzisoxazole-5-carboxylic a...	179	C8H5NO4	1000404-80-3	20
4			trans-.beta.-Ionone	192	C13H20O	000079-77-6	15
5			Octamethyl-1,4-cyclohexadiene	192	C14H24	172684-93-4	15



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
 Data File : BP010038.D
 Acq On : 22 Apr 2022 05:06
 Operator : CG/JU
 Sample : N2490-06
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0J23

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.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, chloro-	5.411	28.9	ng/ul	491421000	1	8.022	340557000	20.0
Benzene, 1,3-di...	7.858	2.5	ng/ul	42791400	1	8.022	340557000	20.0
Benzene, 1,2-di...	8.105	20.0	ng/ul	340557000	1	8.022	340557000	20.0
Benzene, 1,4-di...	8.452	20.5	ng/ul	348679000	1	8.022	340557000	20.0
Benzene, 1-brom...	9.452	8.8	ng/ul	554731	2	10.758	1255270	20.0
m-Chloroaniline	9.781	115.3	ng/ul	7234680	2	10.758	1255270	20.0
Phenol, 2,3-dic...	10.534	38.6	ng/ul	2420060	2	10.758	1255270	20.0
Benzene, 1,2,3-...	10.634	737.1	ng/ul	46262500	2	10.758	1255270	20.0
Benzene, 1,2,4-...	11.146	173.5	ng/ul	10892300	2	10.758	1255270	20.0
Benzenamine, 2,...	13.004	8.6	ng/ul	783258	3	14.569	1830160	20.0
Phenol, 3,4-dic...	13.581	15.3	ng/ul	1403420	3	14.569	1830160	20.0
unknown-01	18.663	6.0	ng/ul	683865	4	17.322	2292910	20.0