

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
 Data File : BP010040.D
 Acq On : 22 Apr 2022 06:13
 Operator : CG/JU
 Sample : N2490-04
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0J21

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: BP010040.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.793	117	120	135	rBV	121952	250526	0.10%	0.027%
2	5.287	358	374	379	rBV2	44235986	128402963	53.00%	13.750%
3	6.622	594	601	611	rVB	327103	561301	0.23%	0.060%
4	7.105	676	683	695	rVB	168734	313296	0.13%	0.034%
5	7.281	704	713	724	rBV	570156	1179945	0.49%	0.126%
6	7.481	739	747	760	rVB	591030	1181389	0.49%	0.127%
7	7.869	797	813	823	rBV	30206329	87355304	36.05%	9.354%
8	8.058	823	845	849	rVV3	44959155	199873382	82.50%	21.403%
9	8.399	880	903	908	rBV2	45894658	242284497	100.00%	25.945%
10	8.652	939	946	956	rVB	532788	864239	0.36%	0.093%
11	9.104	1016	1023	1026	rBV	916893	1502025	0.62%	0.161%
12	9.146	1026	1030	1042	rVB	1467429	2486381	1.03%	0.266%
13	9.446	1071	1081	1090	rBV	686438	1154599	0.48%	0.124%
14	9.757	1129	1134	1140	rVV	480670	904905	0.37%	0.097%
15	9.828	1140	1146	1154	rVV	752509	1357017	0.56%	0.145%
16	10.369	1231	1238	1246	rBV	949190	1772410	0.73%	0.190%
17	10.663	1272	1288	1293	rBV2	41733026	132019558	54.49%	14.137%
18	10.769	1300	1306	1313	rVB	946667	1404726	0.58%	0.150%
19	10.887	1316	1326	1341	rVB2	766138	1627317	0.67%	0.174%
20	11.169	1361	1374	1380	rBV	29412820	64885150	26.78%	6.948%
21	11.346	1397	1404	1420	rVB	1589663	2621988	1.08%	0.281%
22	11.557	1432	1440	1452	rBV	401036	691253	0.29%	0.074%
23	12.034	1512	1521	1530	rBV2	213005	434386	0.18%	0.047%
24	12.775	1629	1647	1654	rBV	2481780	4771034	1.97%	0.511%
25	13.381	1741	1750	1761	rBV	9250927	14233932	5.87%	1.524%
26	13.975	1843	1851	1857	rBV	2137230	3389321	1.40%	0.363%
27	14.028	1857	1860	1869	rVB	199106	306124	0.13%	0.033%
28	14.263	1894	1900	1910	rVB	2380276	3419492	1.41%	0.366%
29	14.569	1945	1952	1963	rBV	1423623	2193525	0.91%	0.235%
30	14.757	1976	1984	1993	rBV	186034	318071	0.13%	0.034%
31	14.892	2001	2007	2014	rBV	184399	261451	0.11%	0.028%
32	15.563	2114	2121	2134	rBV	3057198	4506090	1.86%	0.483%
33	15.675	2134	2140	2149	rVV	1133839	1630058	0.67%	0.175%
34	17.322	2414	2420	2429	rBV	1830551	2586843	1.07%	0.277%
35	17.422	2431	2437	2448	rVB	3412913	4873557	2.01%	0.522%
36	17.569	2458	2462	2468	rBV	191931	294494	0.12%	0.032%

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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

37	18.663	2644	2648	2653	rVB2	307307	398909	0.16%	0.043%
38	19.651	2810	2816	2825	rBV	4426038	5802739	2.40%	0.621%
39	21.104	3060	3063	3068	rBV2	211928	272995	0.11%	0.029%
40	21.404	3108	3114	3119	rBV	1843187	2589026	1.07%	0.277%
41	23.698	3496	3504	3510	rBV	2156234	4526089	1.87%	0.485%
42	23.851	3523	3530	3542	rVB	1101851	2338792	0.97%	0.250%

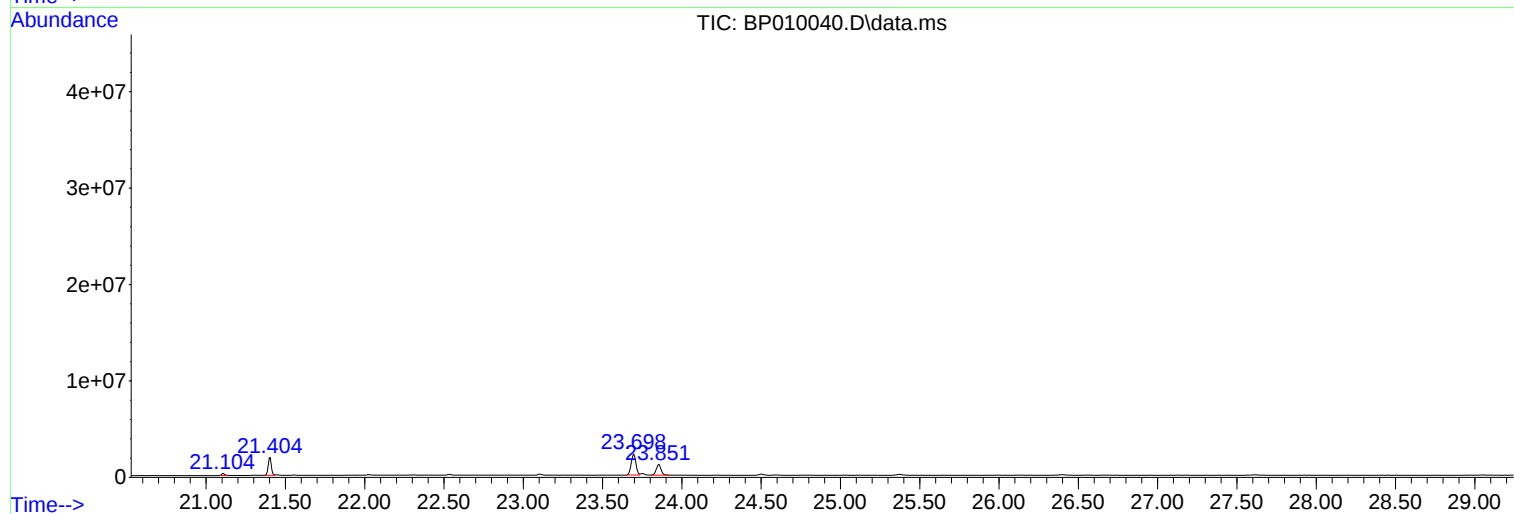
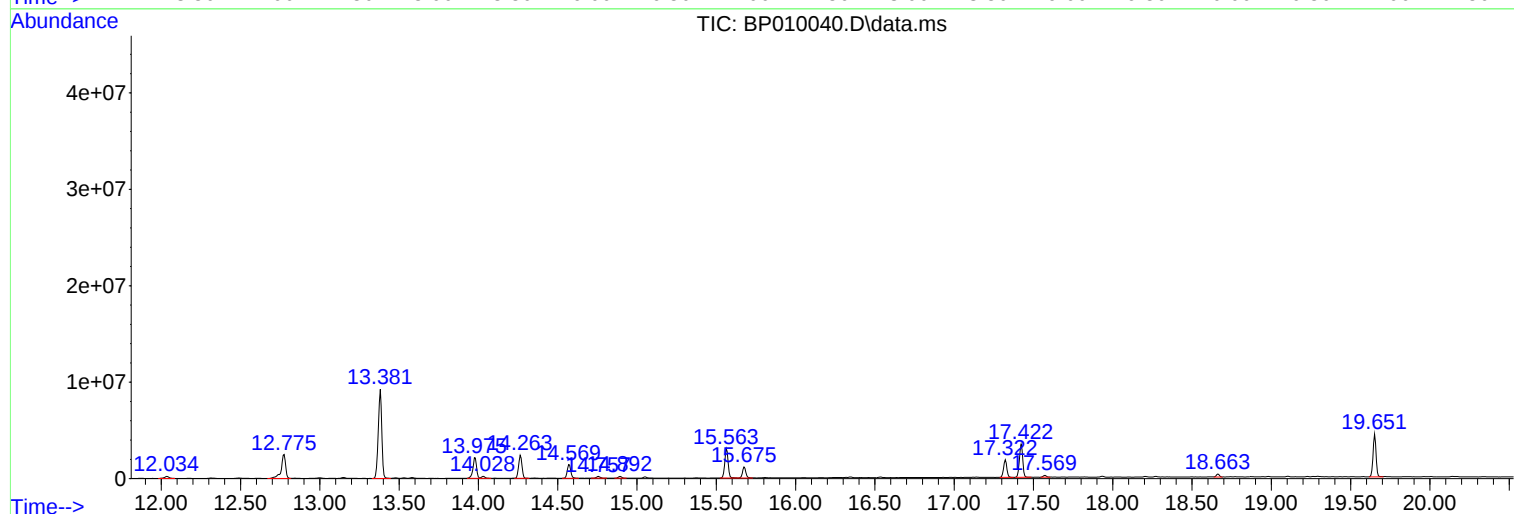
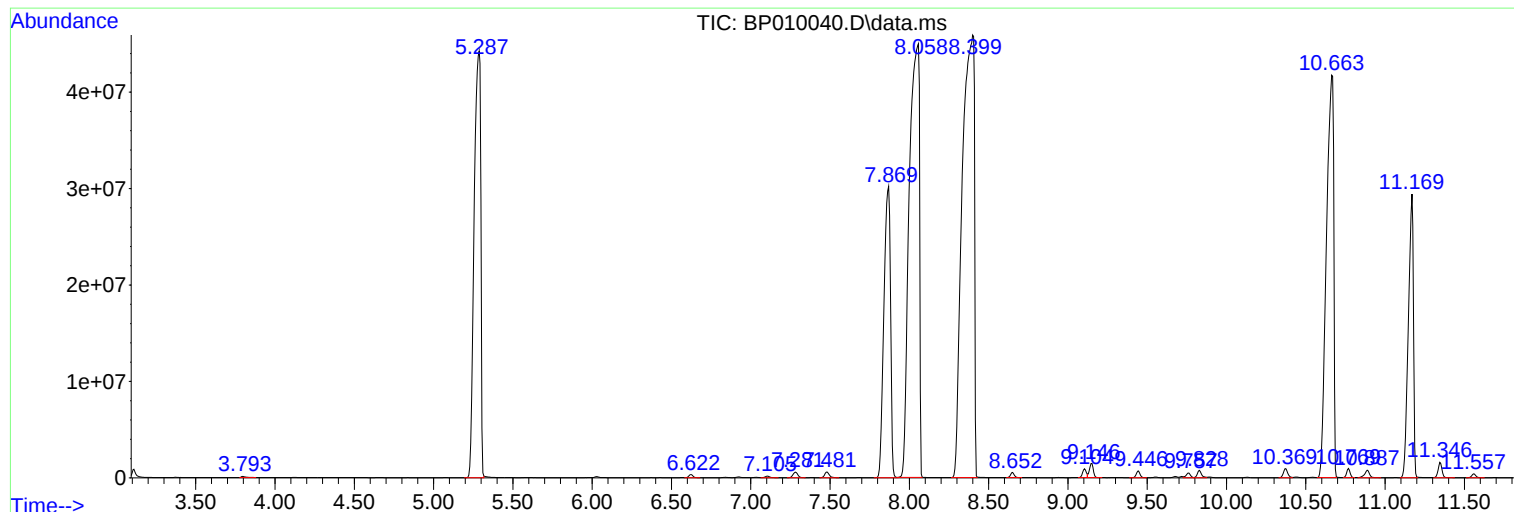
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

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



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Sample : N2490-04
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
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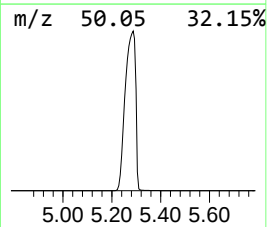
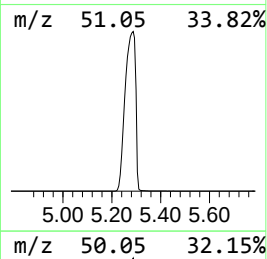
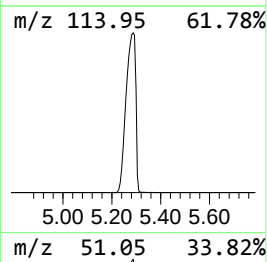
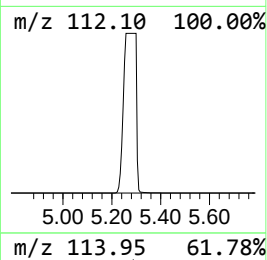
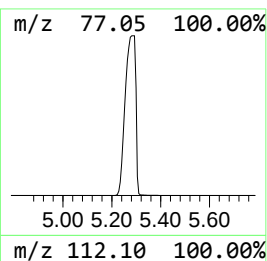
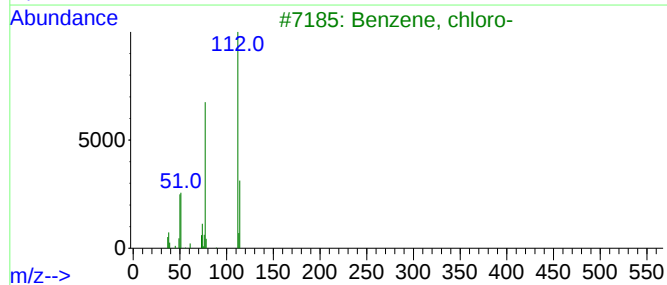
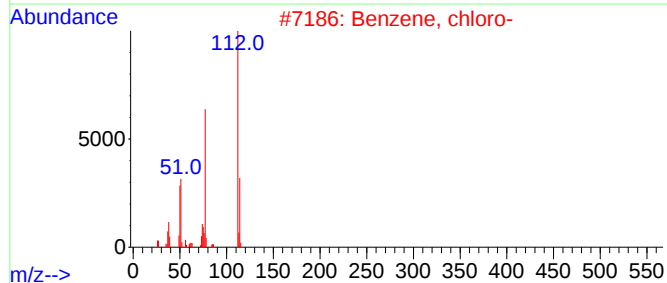
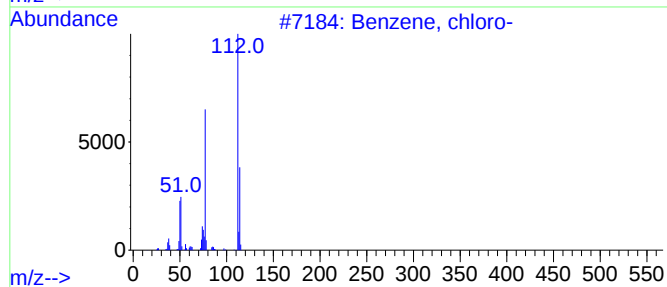
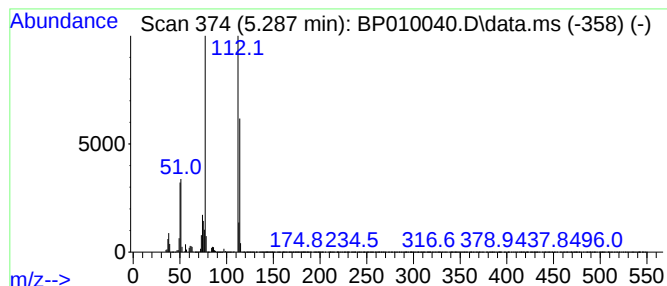
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Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzene, chloro- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.287	12.85 ng/ul	128403000	1,4-Dichlorobenzene-d4	7.981

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, chloro-	112	C6H5Cl	000108-90-7	94
2			Benzene, chloro-	112	C6H5Cl	000108-90-7	93
3			Benzene, chloro-	112	C6H5Cl	000108-90-7	93
4			Benzene, chloro-	112	C6H5Cl	000108-90-7	90
5			Benzene, chloro-	112	C6H5Cl	000108-90-7	90



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Sample : N2490-04
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0J21

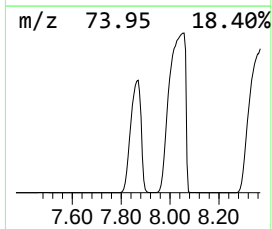
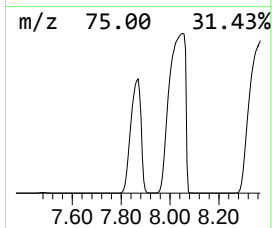
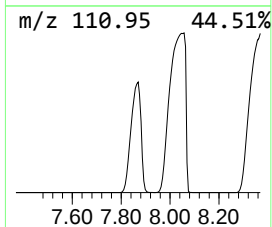
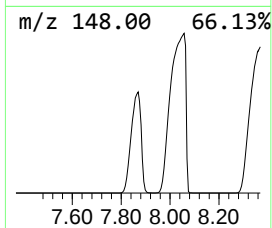
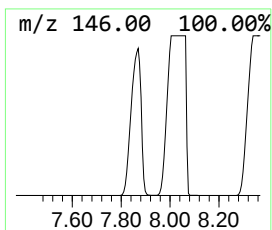
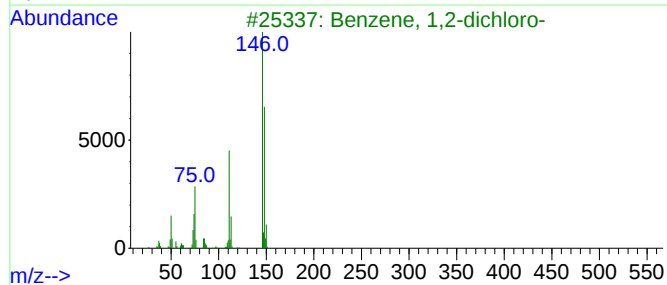
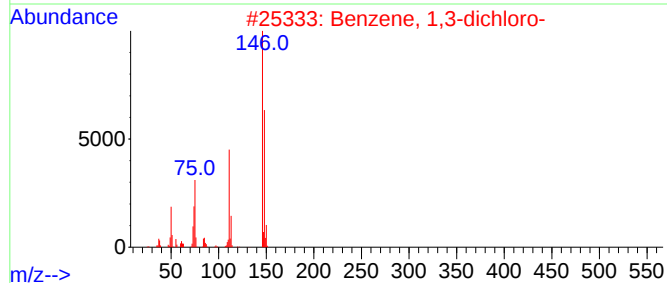
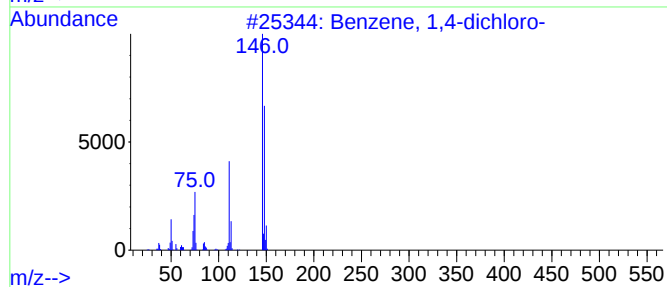
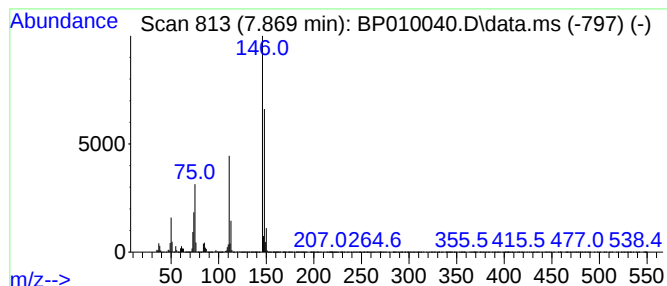
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzene, 1,4-dichloro- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.869	8.74 ng/ul	87355300	1,4-Dichlorobenzene-d4	7.981

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	98
5			Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	97



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Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0J21

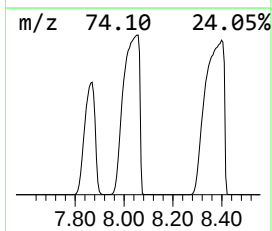
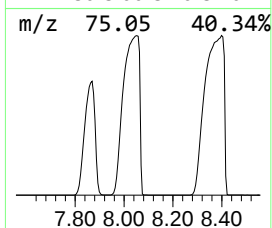
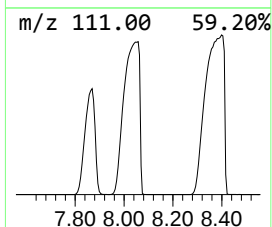
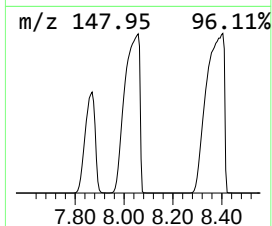
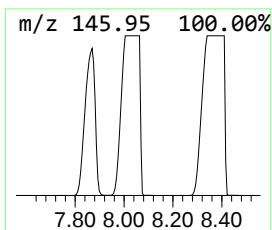
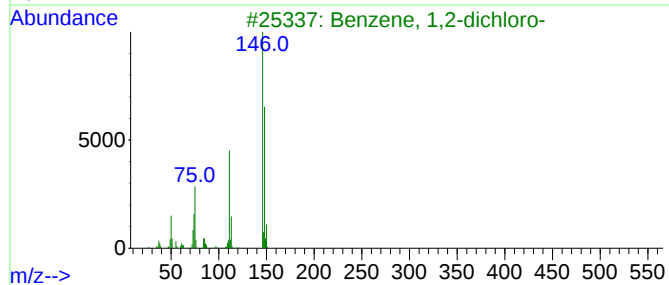
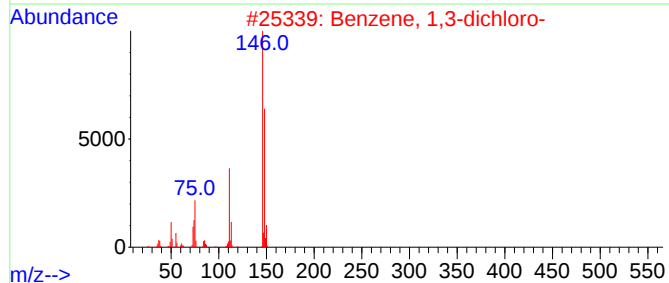
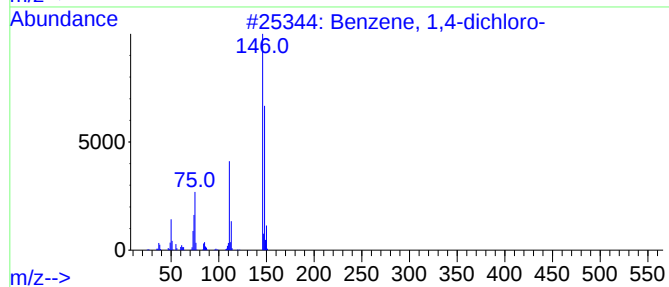
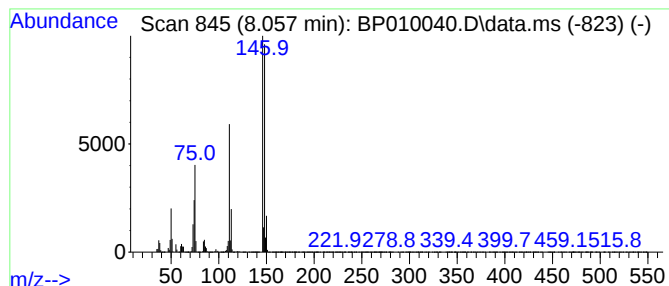
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Benzene, 1,3-dichloro- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.057	20.00 ng/ul	199873000	1,4-Dichlorobenzene-d4	7.981

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	96
3			Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	96
4			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	95
5			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	95



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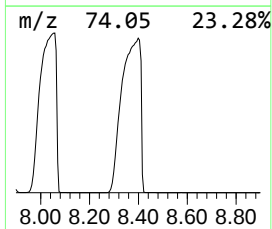
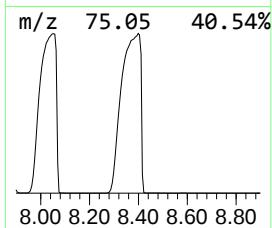
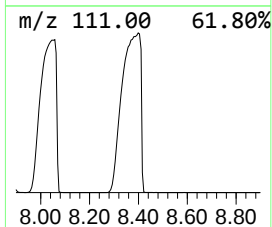
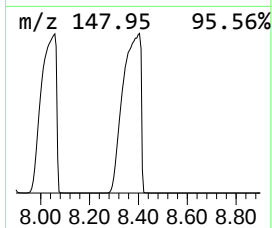
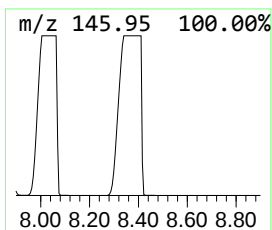
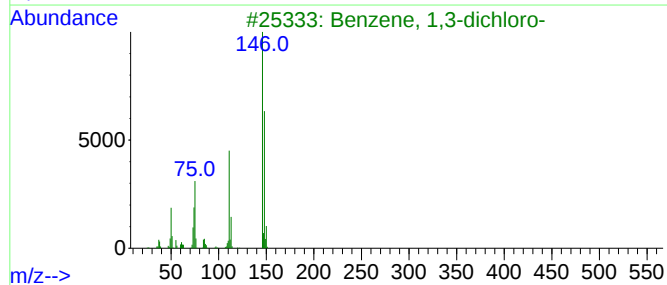
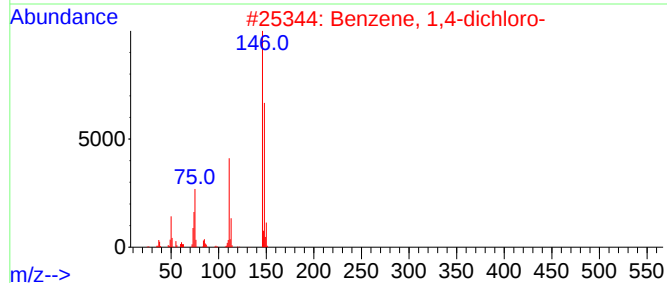
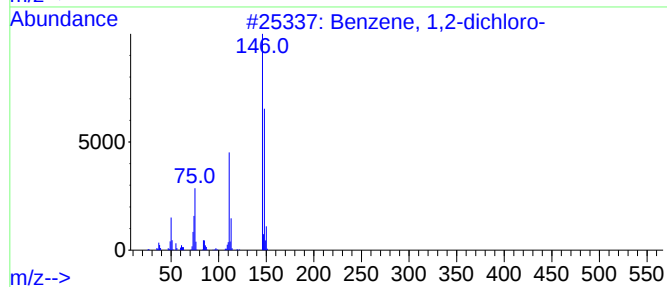
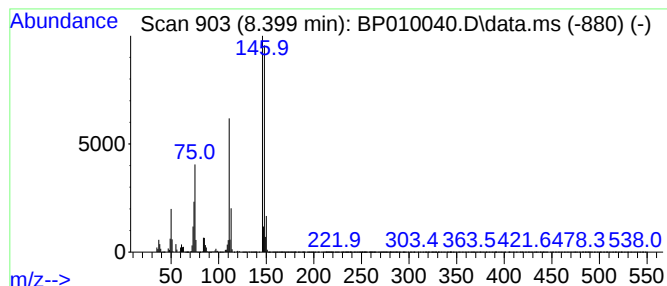
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.399	24.24 ng/ul	242284000	1,4-Dichlorobenzene-d4	7.981

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	96
2			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	96
3			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	95
4			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	93
5			Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	93



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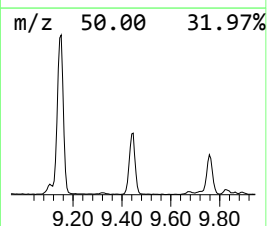
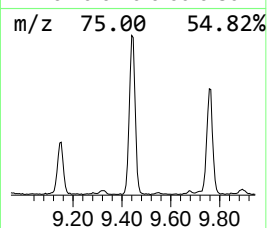
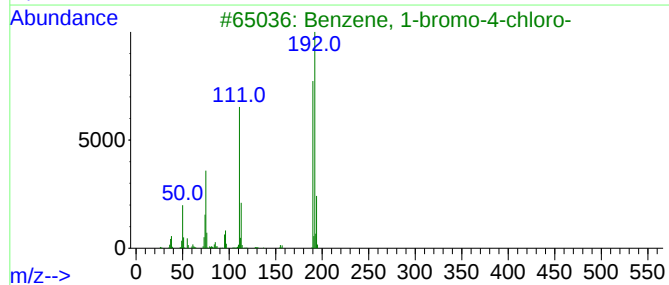
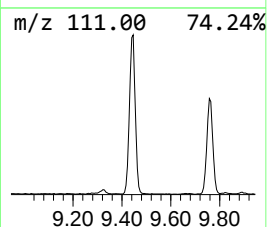
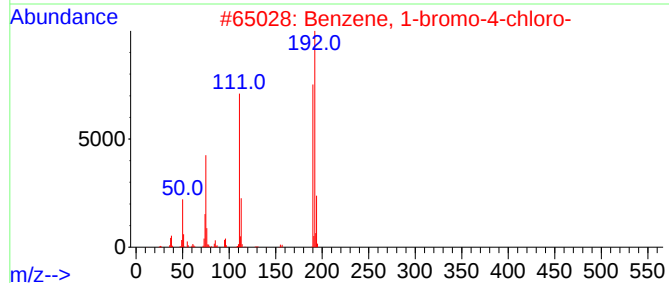
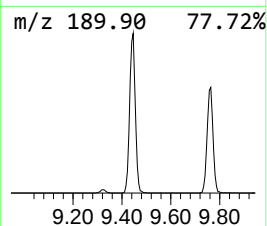
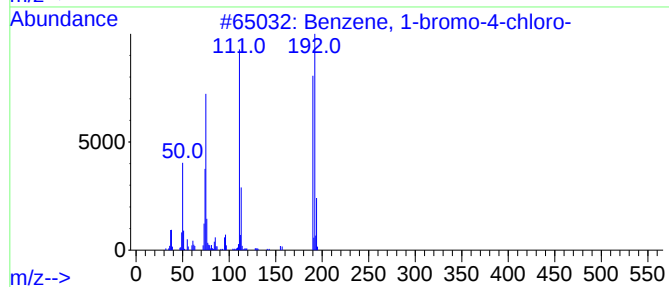
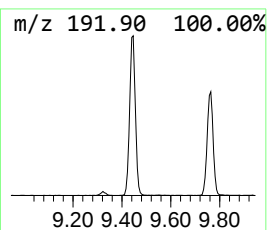
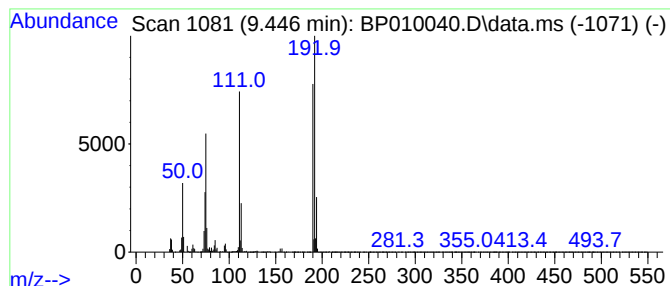
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Benzene, 1-bromo-4-chloro- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.446	16.44 ng/ul	1154600	Naphthalene-d8	10.769

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-4-chloro-	190	C6H4BrCl	000106-39-8	98
3			Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	96
4			Benzene, 1-bromo-3-chloro-	190	C6H4BrCl	000108-37-2	95
5			4-Chlorophenylselenol	192	C6H5ClSe	016645-10-6	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010040.D
Acq On : 22 Apr 2022 06:13
Operator : CG/JU
Sample : N2490-04
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0J21

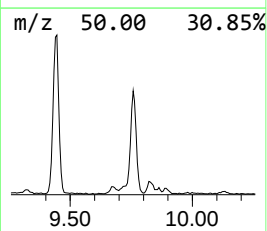
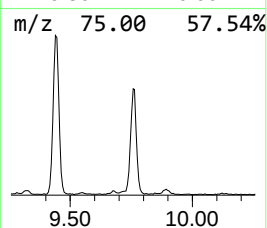
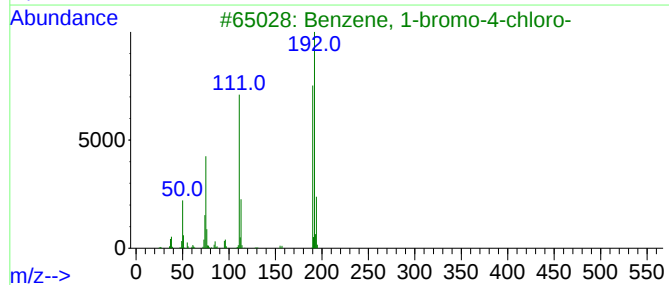
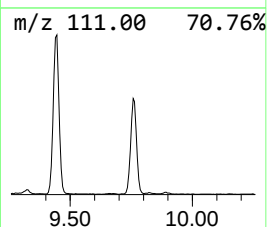
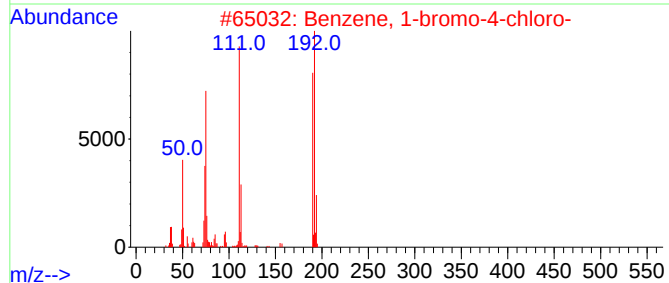
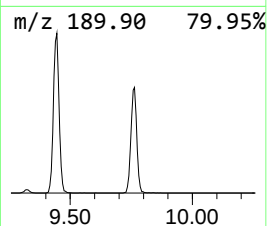
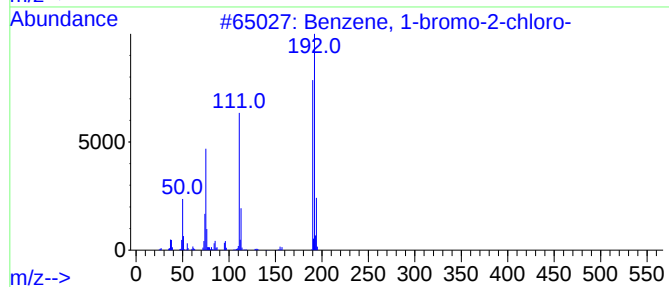
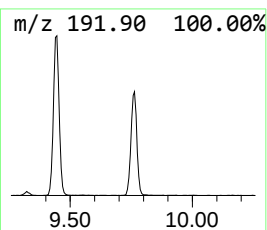
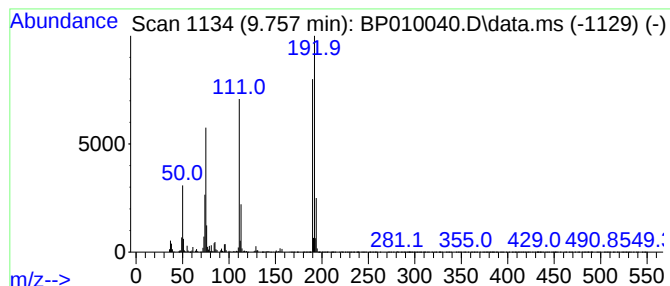
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Quant Title : SVOA CALIBRATION

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.
9.757	12.88 ng/ul	904905	Naphthalene-d8	10.769

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-4-chloro-	190	C6H4BrCl	000106-39-8	96
4			4-Chlorophenylselenol	192	C6H5ClSe	016645-10-6	95
5			Benzene, 1-bromo-3-chloro-	190	C6H4BrCl	000108-37-2	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010040.D
Acq On : 22 Apr 2022 06:13
Operator : CG/JU
Sample : N2490-04
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0J21

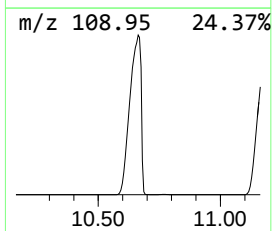
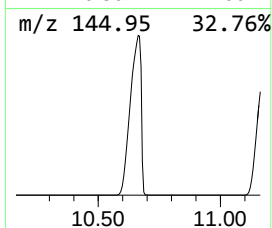
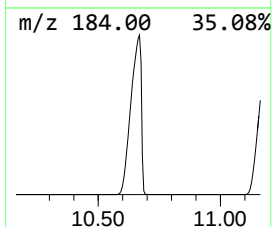
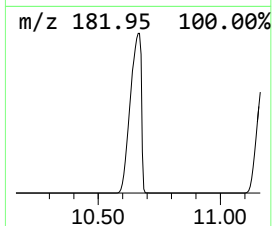
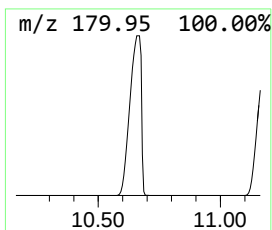
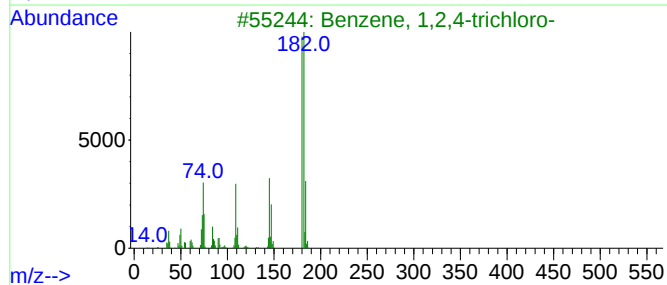
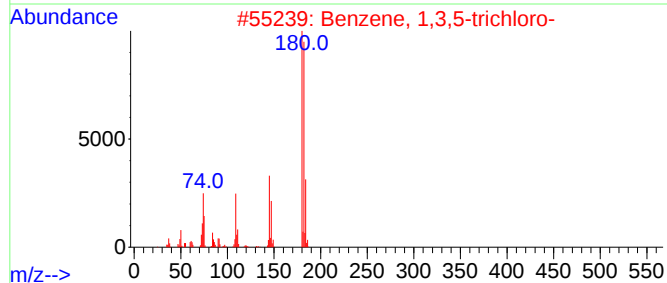
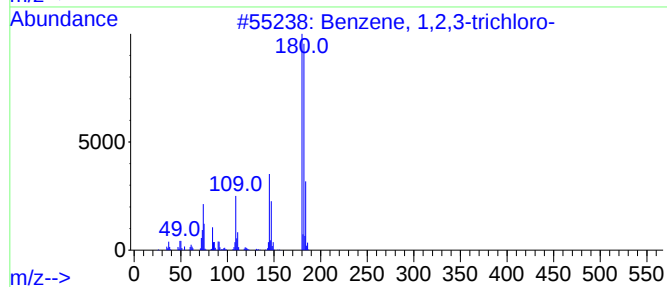
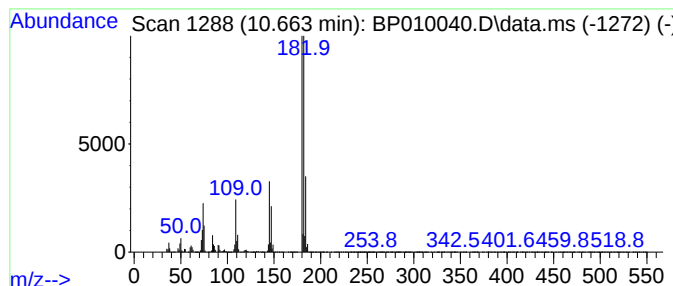
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.663	1879.65 ng/ul	132020000	Naphthalene-d8	10.769

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, 1,2,4-trichloro-	180	C6H3Cl3	000120-82-1	97
4			Benzene, 1,2,3-trichloro-	180	C6H3Cl3	000087-61-6	97
5			Benzene, 1,3,5-trichloro-	180	C6H3Cl3	000108-70-3	97



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010040.D
Acq On : 22 Apr 2022 06:13
Operator : CG/JU
Sample : N2490-04
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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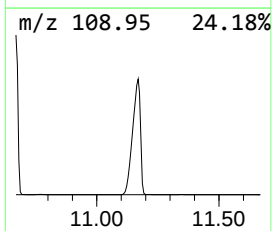
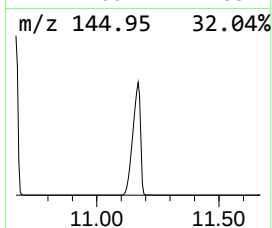
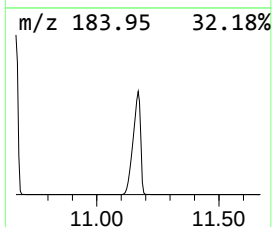
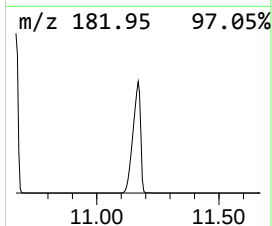
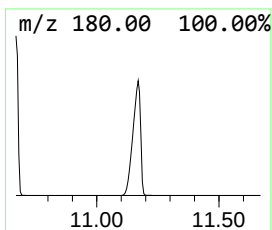
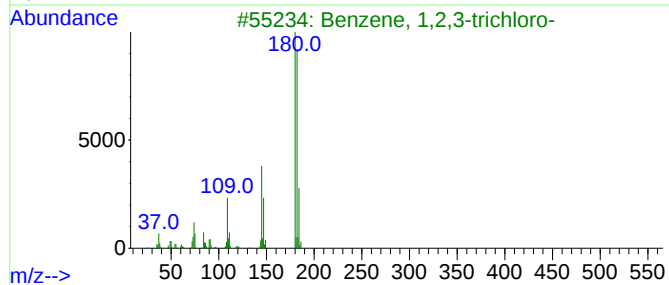
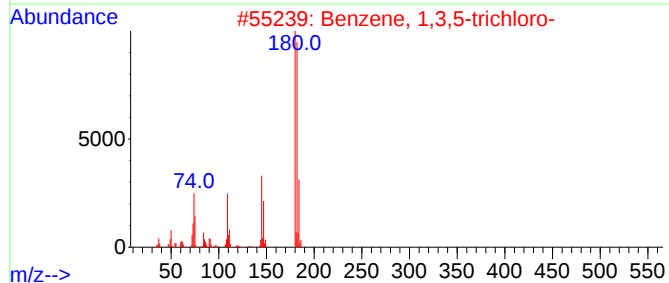
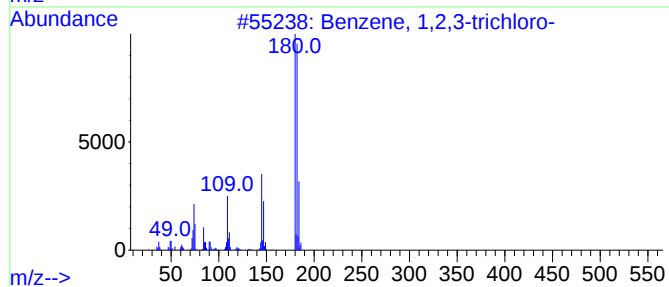
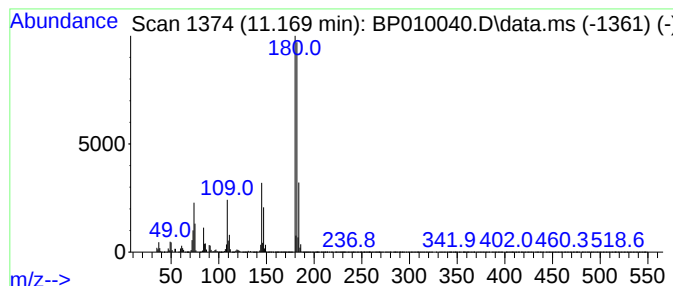
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Benzene, 1,3,5-trichloro- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.169	923.81 ng/ul	64885200	Naphthalene-d8	10.769

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, 1,2,4-trichloro-	180	C6H3Cl3	000120-82-1	97
5			Benzene, 1,2,3-trichloro-	180	C6H3Cl3	000087-61-6	97



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010040.D
Acq On : 22 Apr 2022 06:13
Operator : CG/JU
Sample : N2490-04
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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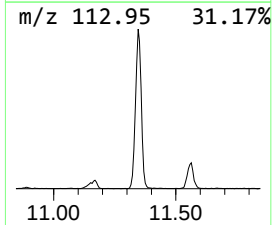
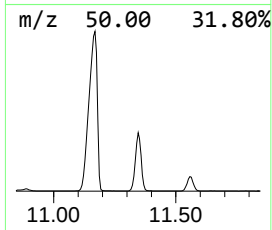
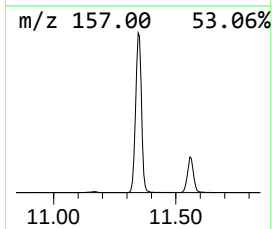
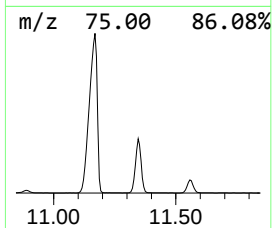
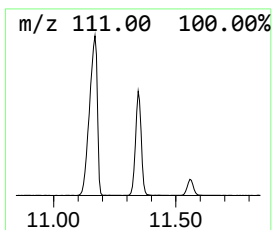
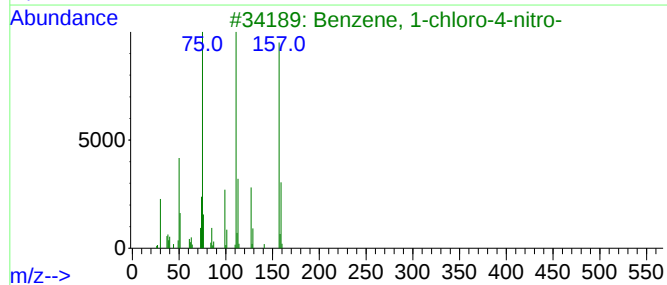
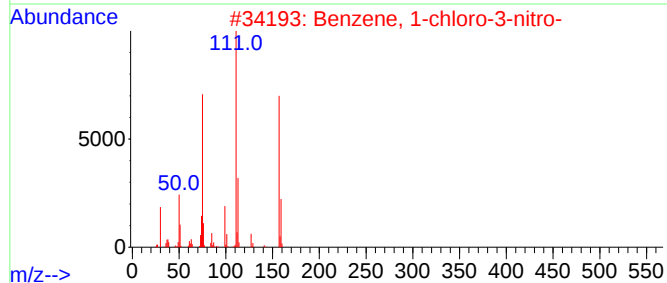
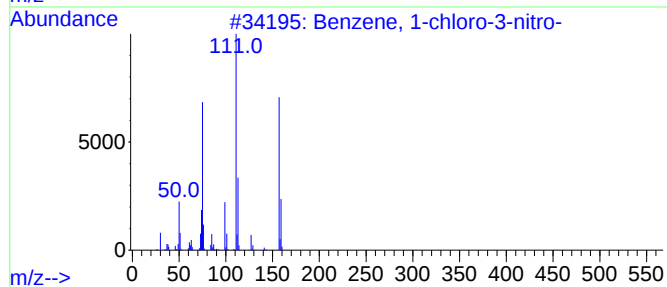
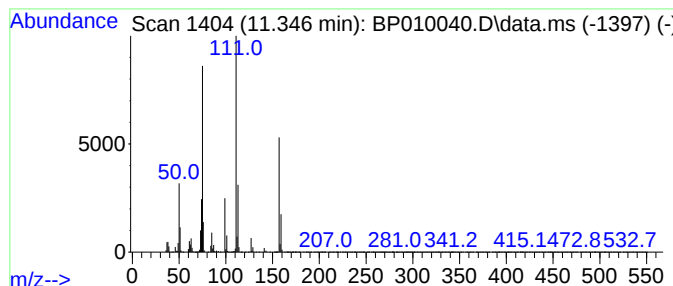
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Benzene, 1-chloro-3-nitro- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.346	37.33 ng/ul	2621990	Naphthalene-d8	10.769

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-3-nitro-	157	C6H4ClNO2	000121-73-3	94
3			Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	94
4			Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	94
5			Benzene, 1-chloro-3-nitro-	157	C6H4ClNO2	000121-73-3	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010040.D
Acq On : 22 Apr 2022 06:13
Operator : CG/JU
Sample : N2490-04
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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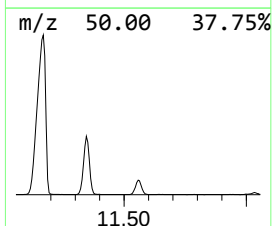
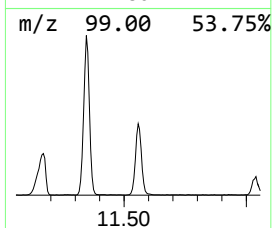
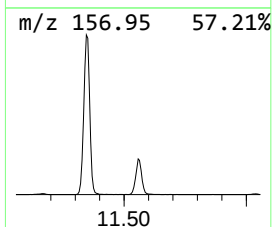
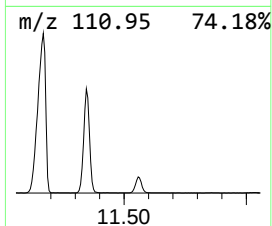
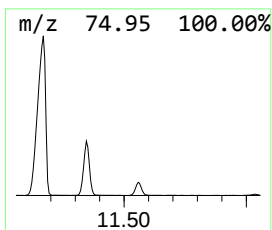
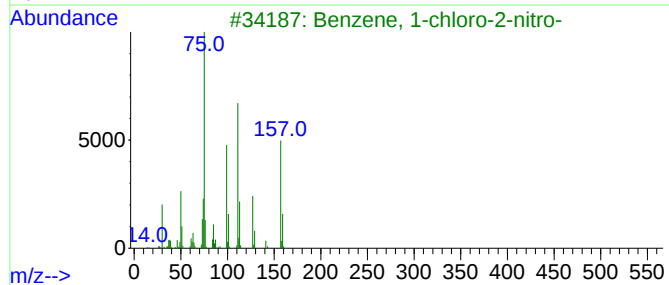
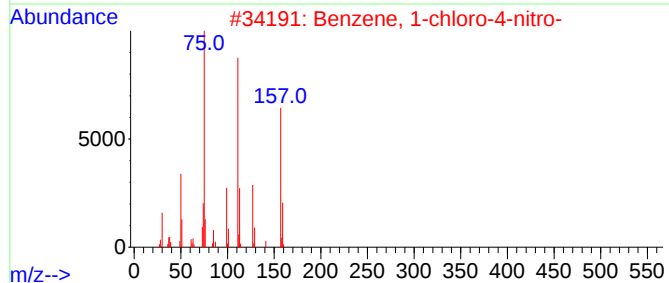
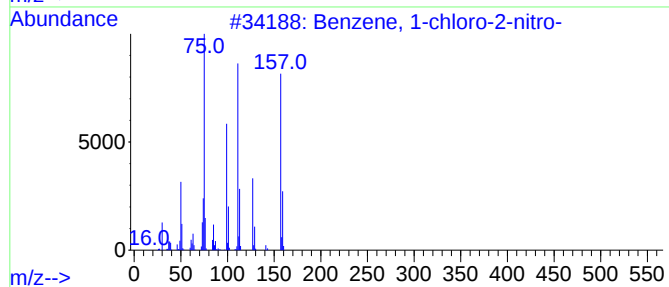
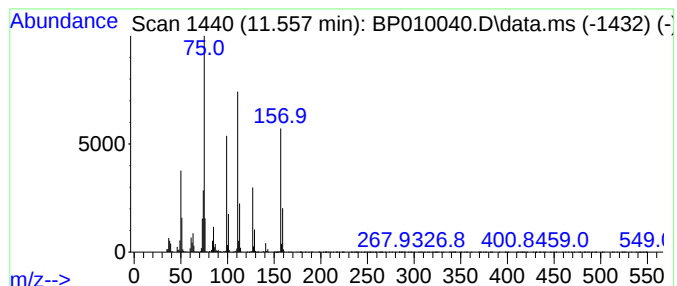
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Benzene, 1-chloro-2-nitro- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.557	9.84 ng/ul	691253	Naphthalene-d8	10.769

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-chloro-2-nitro-	157	C6H4ClNO2	000088-73-3	97
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	95
5			Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010040.D
Acq On : 22 Apr 2022 06:13
Operator : CG/JU
Sample : N2490-04
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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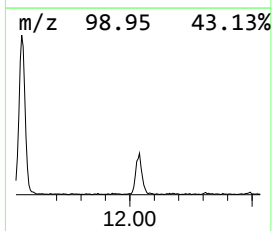
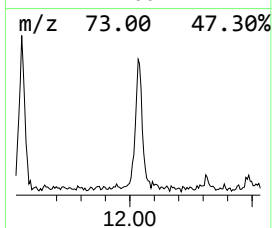
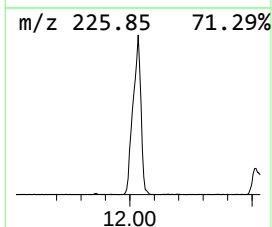
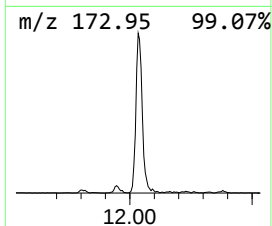
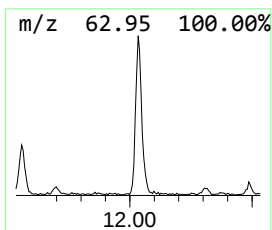
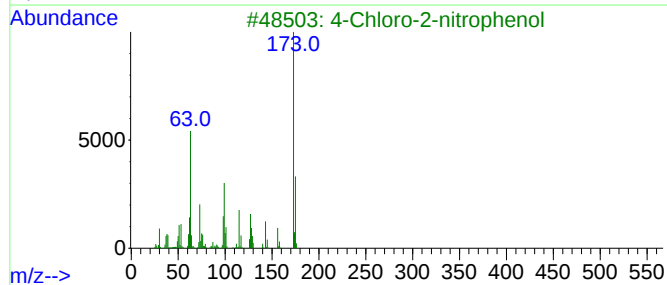
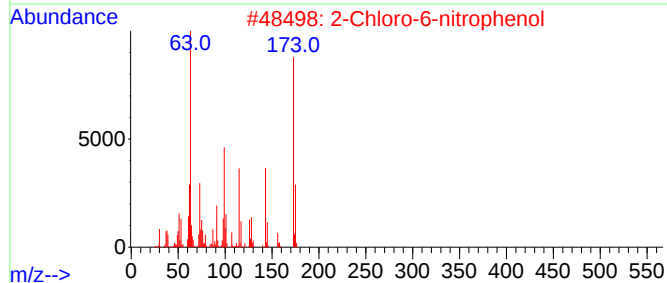
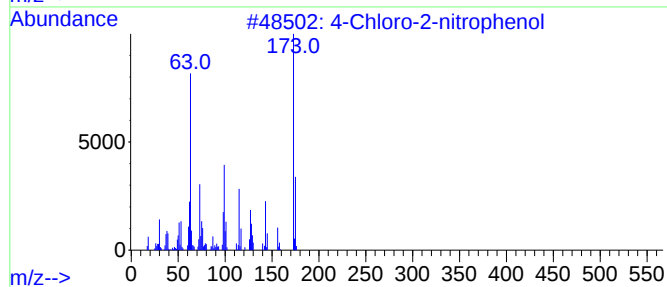
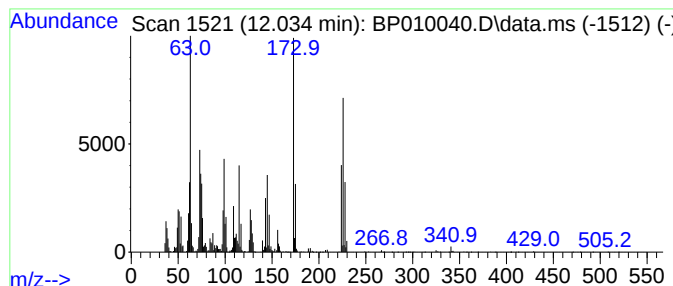
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 4-Chloro-2-nitrophenol Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.034	6.18 ng/ul	434386	Naphthalene-d8	10.769

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-Chloro-2-nitrophenol	173	C6H4ClNO3	000089-64-5	99
2			2-Chloro-6-nitrophenol	173	C6H4ClNO3	000603-86-1	93
3			4-Chloro-2-nitrophenol	173	C6H4ClNO3	000089-64-5	93
4			Phenol, 5-chloro-2-nitro-	173	C6H4ClNO3	000611-07-4	90
5			2-Chloro-5-nitrophenol	173	C6H4ClNO3	000619-10-3	89



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010040.D
Acq On : 22 Apr 2022 06:13
Operator : CG/JU
Sample : N2490-04
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0J21

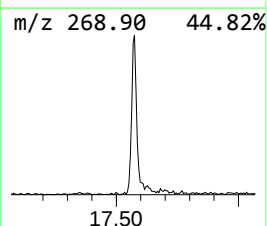
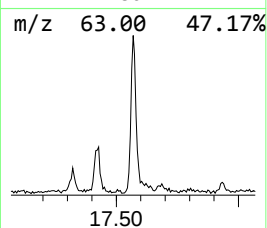
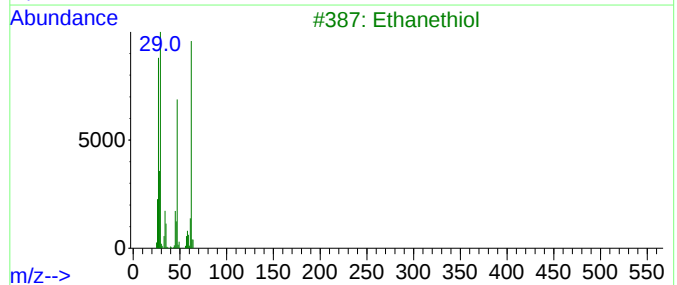
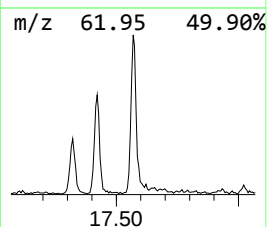
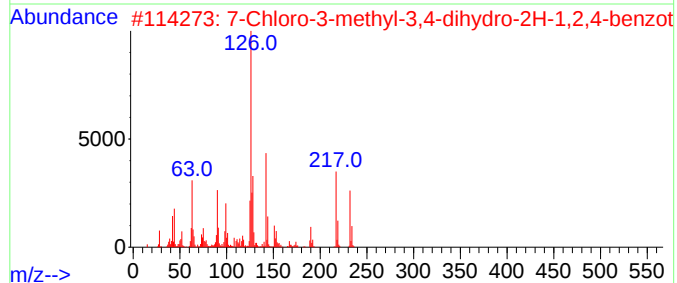
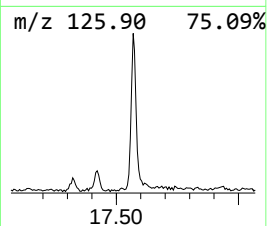
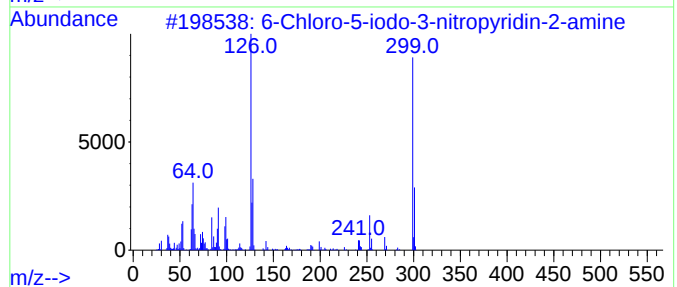
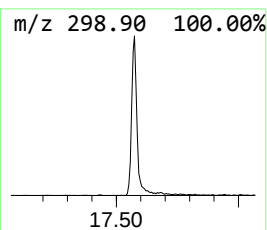
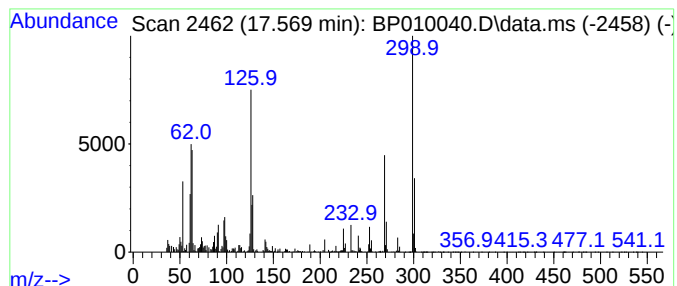
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 unknown-01 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.569	2.28 ng/ul	294494	Phenanthrene-d10	17.322

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			6-Chloro-5-iodo-3-nitropyridin-2...	299	C5H3ClIN3O2	790692-90-9	38
2			7-Chloro-3-methyl-3,4-dihydro-2H...	232	C8H9ClN2O2S	022503-72-6	17
3			Ethanethiol	62	C2H6S	000075-08-1	9
4			Ethanethiol	62	C2H6S	000075-08-1	9
5			Ethene, chloro-	62	C2H3Cl	000075-01-4	9



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010040.D
Acq On : 22 Apr 2022 06:13
Operator : CG/JU
Sample : N2490-04
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0J21

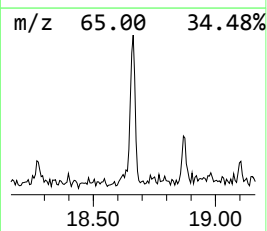
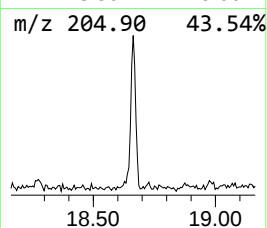
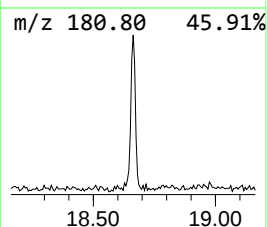
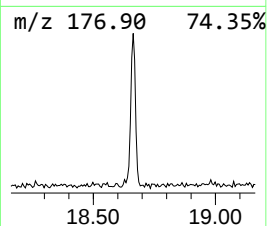
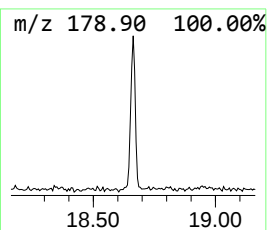
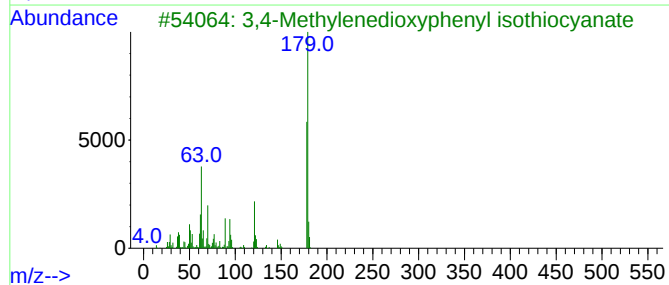
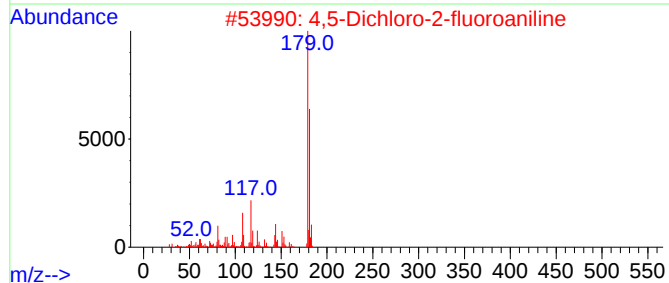
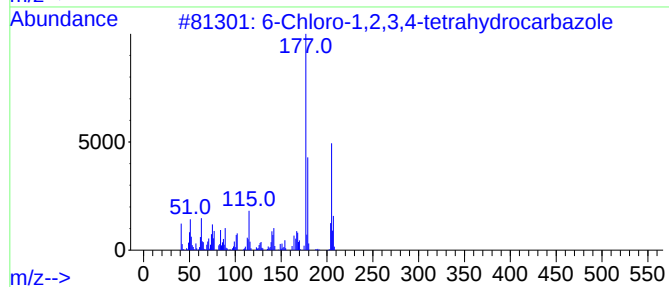
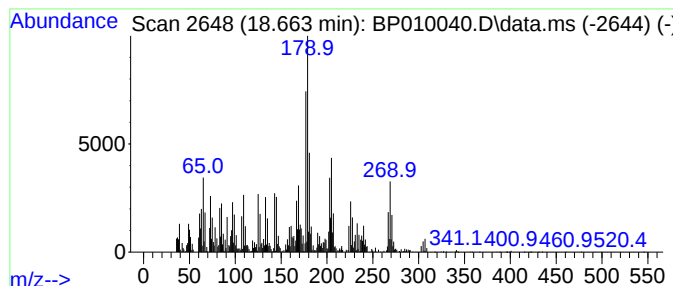
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 unknown-02 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.663	3.08 ng/ul	398909	Phenanthrene-d10	17.322

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			6-Chloro-1,2,3,4-tetrahydrocarba...	205	C12H12ClN	036684-65-8	42
2			4,5-Dichloro-2-fluoroaniline	179	C6H4Cl2FN	002729-36-4	25
3			3,4-Methylenedioxyphenyl isothio...	179	C8H5NO2S	113504-93-1	25
4			2,1-benzisoxazole-5-carboxylic a...	179	C8H5NO4	1000404-80-3	25
5			2,5-Dichloro-4-fluoroaniline	179	C6H4Cl2FN	002729-37-5	22



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010040.D
Acq On : 22 Apr 2022 06:13
Operator : CG/JU
Sample : N2490-04
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0J21

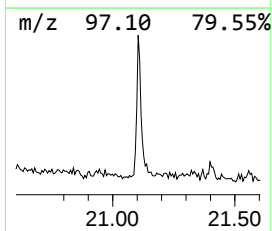
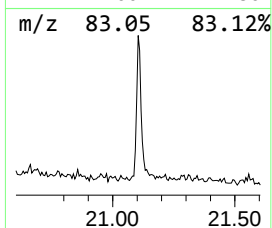
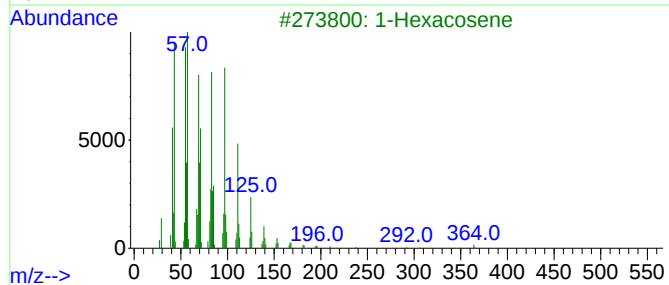
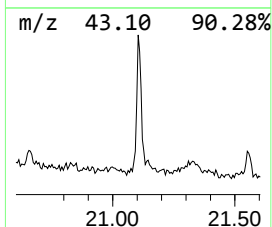
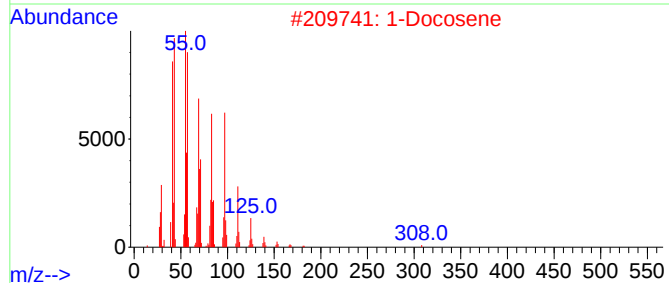
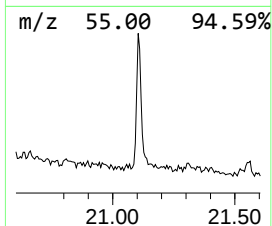
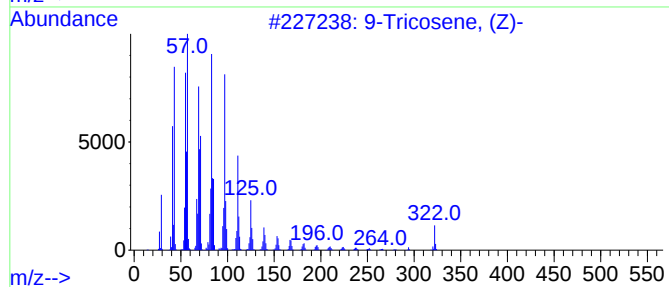
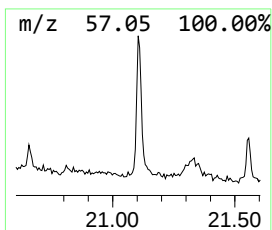
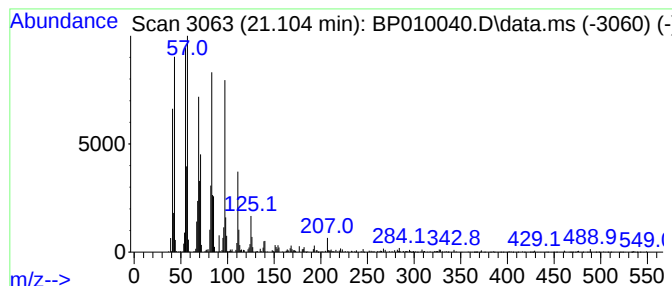
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

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

Peak Number 14 (DEL) Alkane: Straight-Chai... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.104	2.11 ng/ul	272995	Chrysene-d12	21.404

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Tricosene, (Z)-	322	C23H46	027519-02-4	97	
2	1-Docosene	308	C22H44	001599-67-3	96	
3	1-Hexacosene	364	C26H52	018835-33-1	94	
4	1-Tricosene	322	C23H46	018835-32-0	91	
5	1-Tetracosene	336	C24H48	010192-32-2	91	



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
 Data File : BP010040.D
 Acq On : 22 Apr 2022 06:13
 Operator : CG/JU
 Sample : N2490-04
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0J21

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-	5.287	12.9	ng/ul	128403000	1	7.981	199873000	20.0
Benzene, 1,4-di...	7.869	8.7	ng/ul	87355300	1	7.981	199873000	20.0
Benzene, 1,3-di...	8.057	20.0	ng/ul	199873000	1	7.981	199873000	20.0
Benzene, 1,2-di...	8.399	24.2	ng/ul	242284000	1	7.981	199873000	20.0
Benzene, 1-brom...	9.446	16.4	ng/ul	1154600	2	10.769	1404730	20.0
Benzene, 1-brom...	9.757	12.9	ng/ul	904905	2	10.769	1404730	20.0
Benzene, 1,2,3-...	10.663	1879.7	ng/ul	132020000	2	10.769	1404730	20.0
Benzene, 1,3,5-...	11.169	923.8	ng/ul	64885200	2	10.769	1404730	20.0
Benzene, 1-chlo...	11.346	37.3	ng/ul	2621990	2	10.769	1404730	20.0
Benzene, 1-chlo...	11.557	9.8	ng/ul	691253	2	10.769	1404730	20.0
4-Chloro-2-nitr...	12.034	6.2	ng/ul	434386	2	10.769	1404730	20.0
unknown-01	17.569	2.3	ng/ul	294494	4	17.322	2586840	20.0
unknown-02	18.663	3.1	ng/ul	398909	4	17.322	2586840	20.0
(DEL) Alkane: S...	21.104	2.1	ng/ul	272995	5	21.404	2589030	20.0