

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
 Data File : BP010067.D
 Acq On : 22 Apr 2022 22:14
 Operator : CG/JU
 Sample : N2490-07
 Misc :
 ALS Vial : 63 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0J44

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: BP010067.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.364	360	387	391	rBV3	50357432	343990822	100.00%	37.217%
2	6.622	594	601	608	rBV	225357	359830	0.10%	0.039%
3	7.281	705	713	726	rBV	553687	1080787	0.31%	0.117%
4	7.481	738	747	756	rBV	577682	1082235	0.31%	0.117%
5	7.675	772	780	791	rVB	341087	779224	0.23%	0.084%
6	7.852	798	810	822	rBV	9085781	24515961	7.13%	2.652%
7	8.063	822	846	850	rBV3	45811281	236567639	68.77%	25.595%
8	8.387	880	901	905	rBV2	43940025	190794541	55.47%	20.643%
9	8.652	939	946	957	rVB	504058	844555	0.25%	0.091%
10	9.110	1016	1024	1026	rBV	737927	1370124	0.40%	0.148%
11	9.163	1026	1033	1041	rVB	7044568	12776533	3.71%	1.382%
12	9.558	1093	1100	1109	rVB2	297911	526234	0.15%	0.057%
13	9.828	1139	1146	1154	rBV	599017	1007783	0.29%	0.109%
14	10.369	1229	1238	1252	rBV	949765	1670370	0.49%	0.181%
15	10.634	1270	1283	1297	rBV	26073197	53307956	15.50%	5.768%
16	10.757	1297	1304	1309	rVV	850328	1418176	0.41%	0.153%
17	10.881	1317	1325	1344	rVB	546688	1060198	0.31%	0.115%
18	11.134	1358	1368	1384	rBV	2800659	4725285	1.37%	0.511%
19	11.346	1398	1404	1410	rVV	1110736	1935044	0.56%	0.209%
20	11.557	1433	1440	1450	rBV	356818	616887	0.18%	0.067%
21	12.769	1635	1646	1655	rBV	1855945	3106386	0.90%	0.336%
22	13.375	1742	1749	1758	rBV2	3132445	4975708	1.45%	0.538%
23	13.581	1778	1784	1791	rVB	316487	493183	0.14%	0.053%
24	13.975	1844	1851	1857	rBV	2068309	3023061	0.88%	0.327%
25	14.263	1890	1900	1907	rBV	2308394	3425908	1.00%	0.371%
26	14.569	1943	1952	1969	rVB	1413994	2184735	0.64%	0.236%
27	15.563	2114	2121	2135	rBV	3077942	4635413	1.35%	0.502%
28	15.681	2135	2141	2152	rVB	446426	686707	0.20%	0.074%
29	17.322	2412	2420	2429	rBV	1753914	2602704	0.76%	0.282%
30	17.422	2430	2437	2448	rVV	3275026	4847748	1.41%	0.524%
31	17.992	2529	2534	2545	rVB	346166	432848	0.13%	0.047%
32	19.122	2718	2726	2730	rBV	420789	517234	0.15%	0.056%
33	19.245	2739	2747	2752	rBV2	328812	456454	0.13%	0.049%
34	19.651	2810	2816	2833	rBV	3994702	5353514	1.56%	0.579%
35	21.404	3109	3114	3125	rBV	1447682	2029897	0.59%	0.220%
36	23.698	3497	3504	3511	rBV2	1496998	3294416	0.96%	0.356%

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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	23.857	3524	3531	3545	rVB	803770	1783197	0.52%	0.193%
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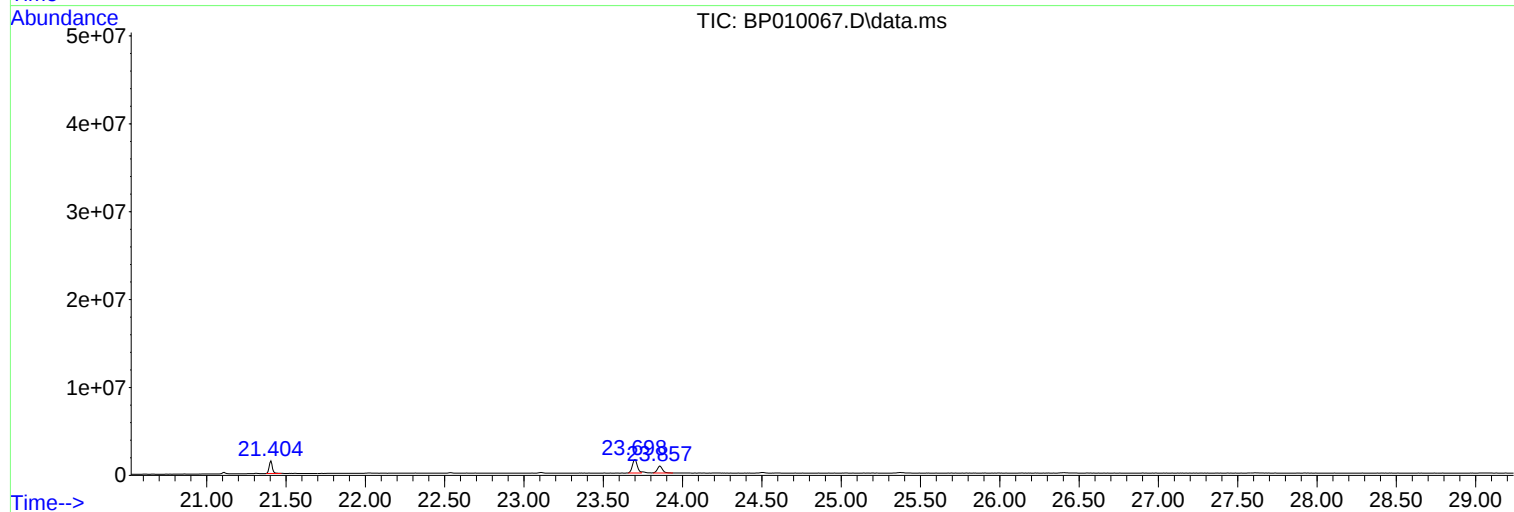
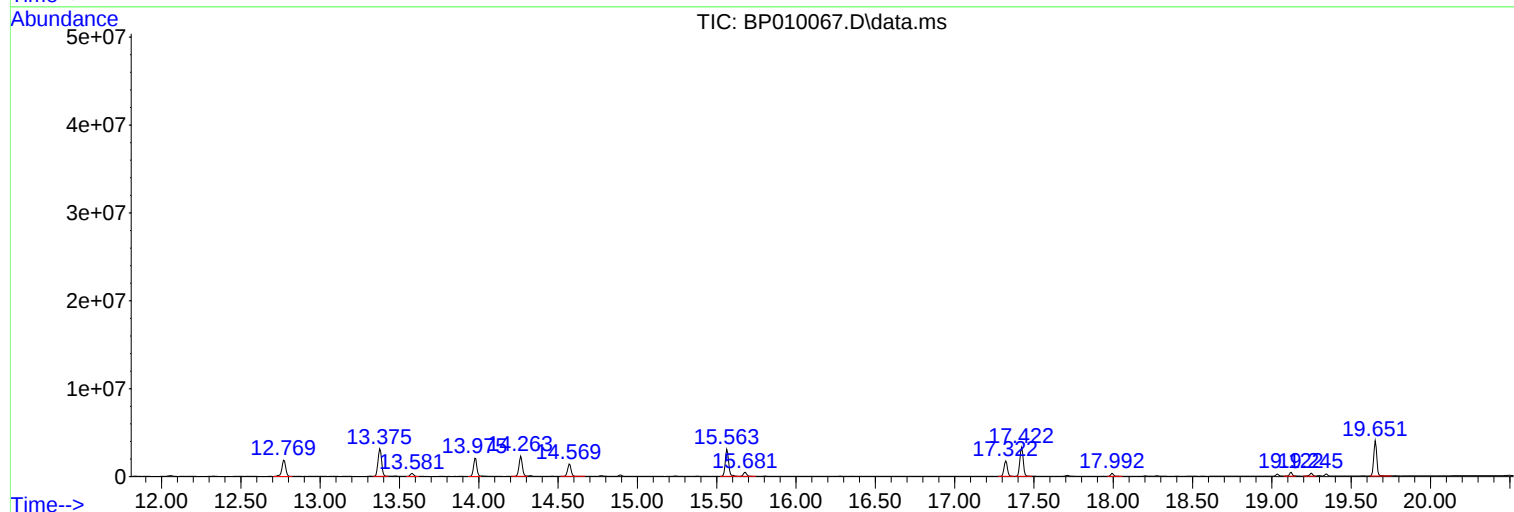
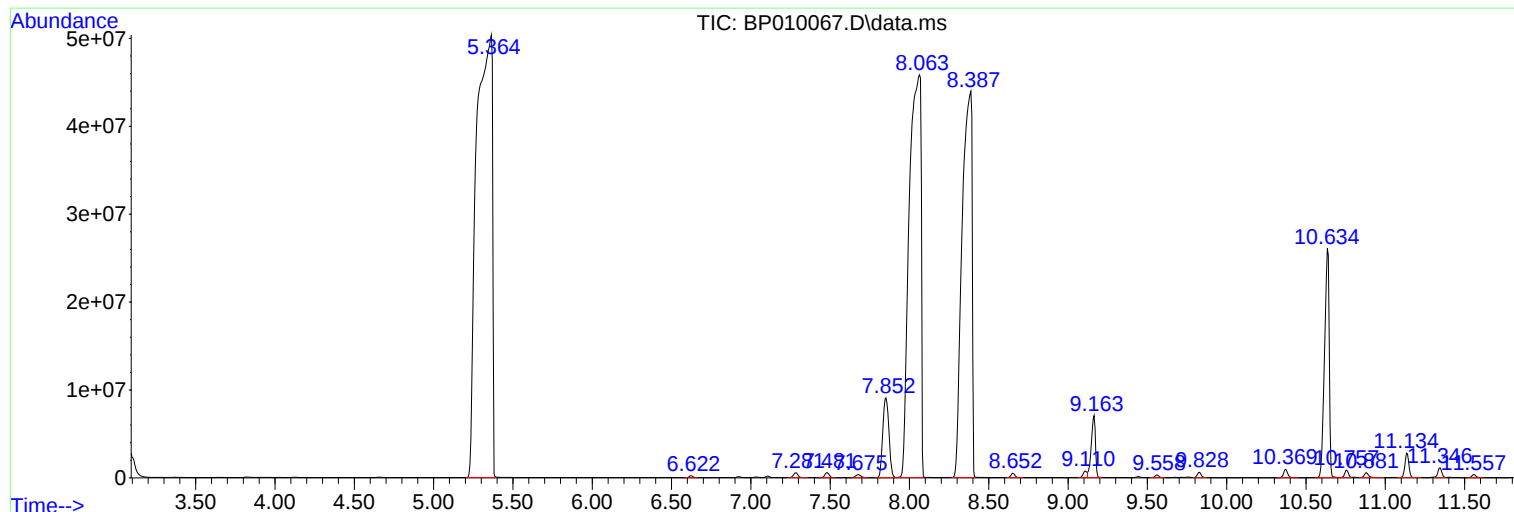
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

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



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Operator : CG/JU
Sample : N2490-07
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Instrument :
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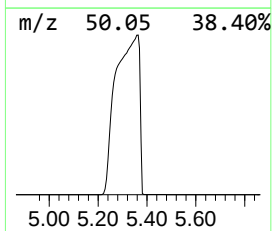
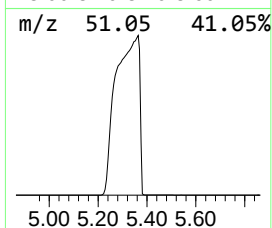
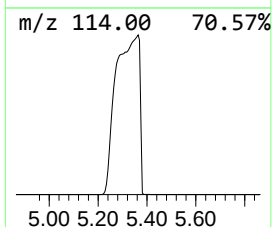
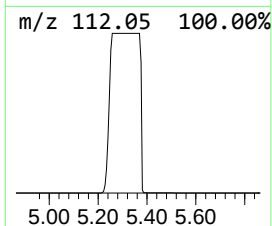
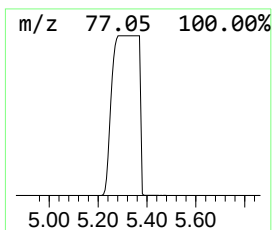
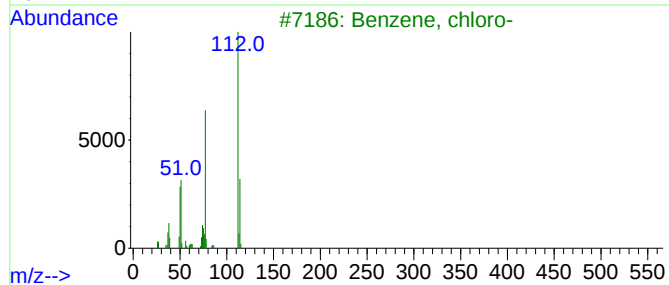
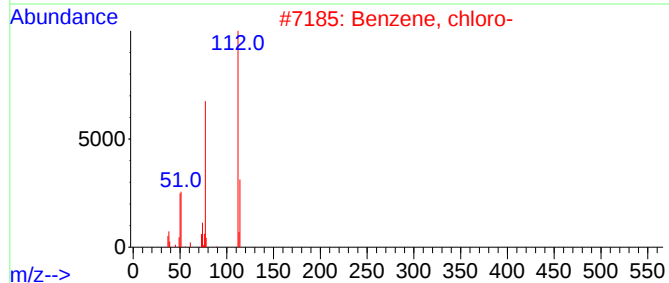
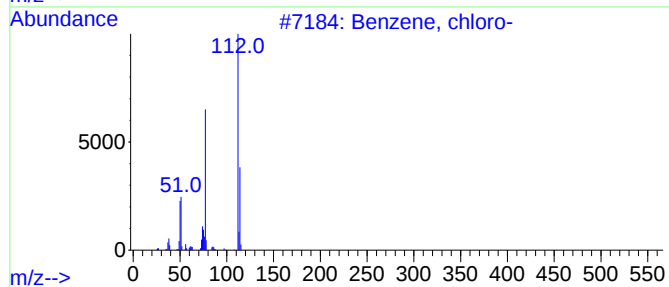
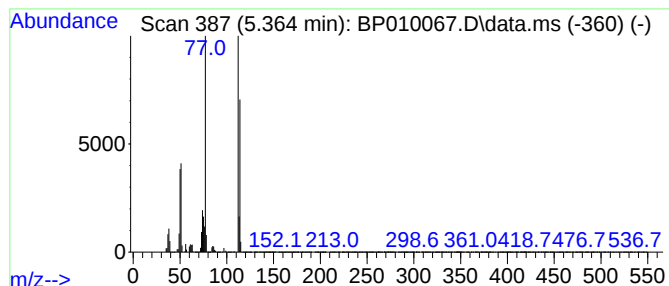
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzene, chloro- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.364	29.08 ng/ul	343991000	1,4-Dichlorobenzene-d4	7.993

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	94
4			Benzene, chloro-	112	C6H5Cl	000108-90-7	78
5			Benzene, chloro-	112	C6H5Cl	000108-90-7	78



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ALS Vial : 63 Sample Multiplier: 1

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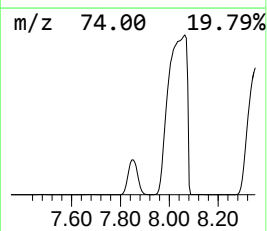
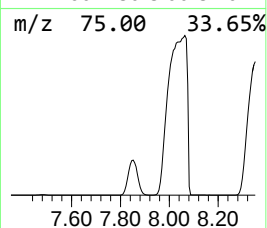
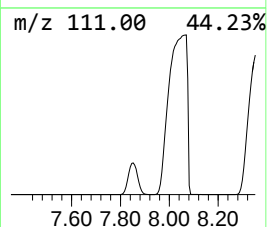
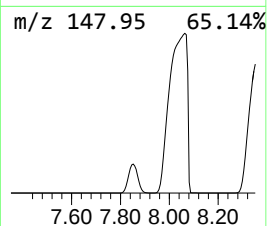
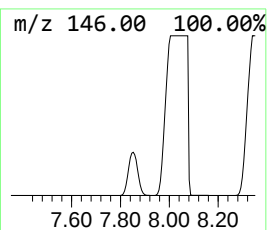
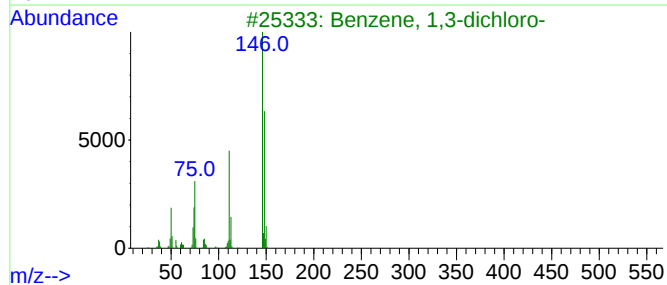
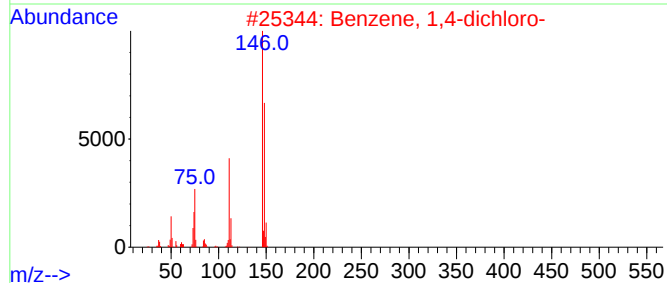
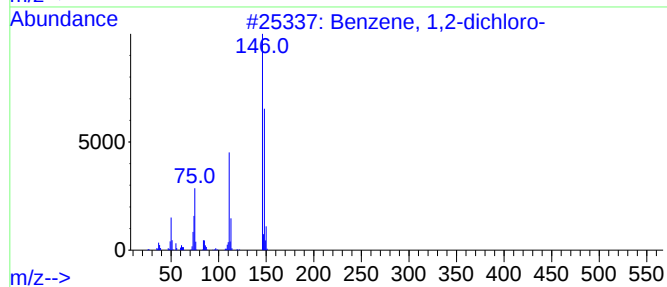
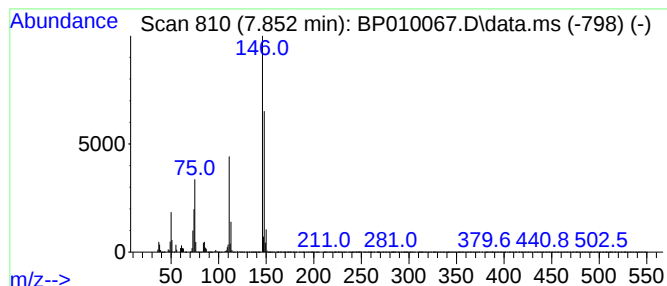
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TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Benzene, 1,2-dichloro- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.852	2.07 ng/ul	24516000	1,4-Dichlorobenzene-d4	7.993

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	98
2			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	98
3			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	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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ALS Vial : 63 Sample Multiplier: 1

Instrument :
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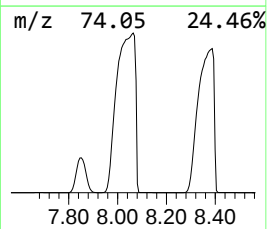
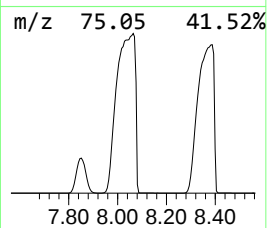
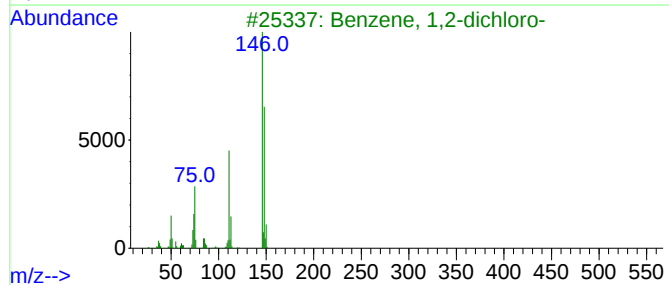
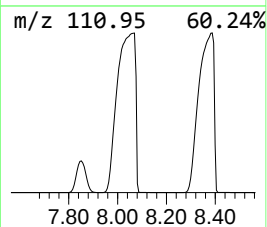
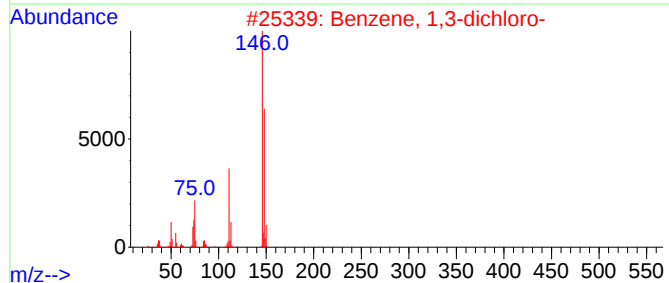
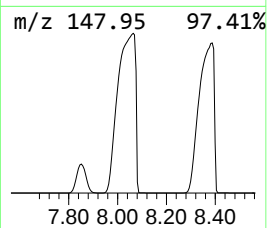
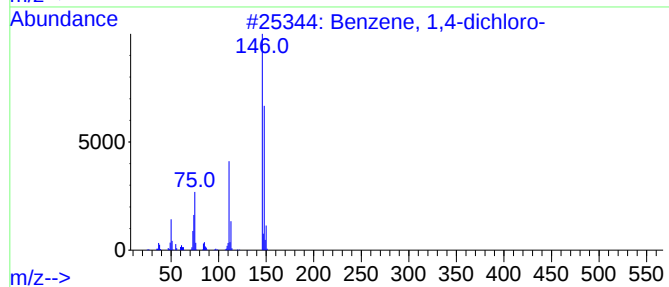
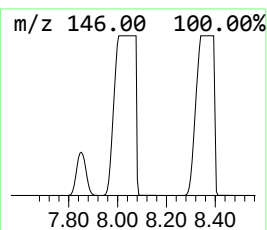
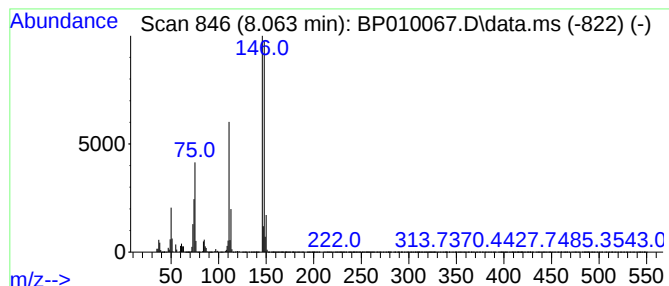
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Benzene, 1,4-dichloro- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.063	20.00 ng/ul	236568000	1,4-Dichlorobenzene-d4	7.993

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	96
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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ClientSampleId :
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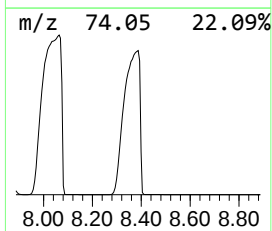
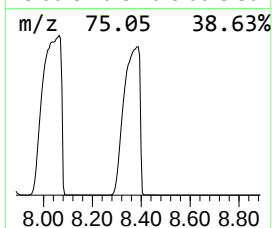
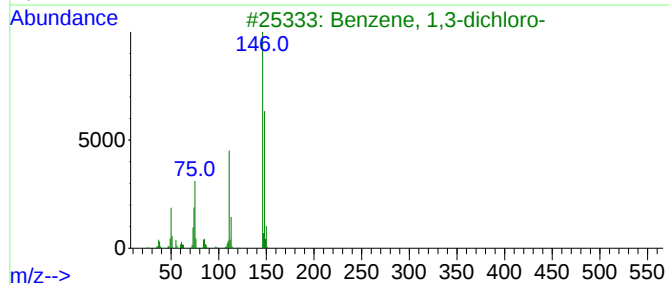
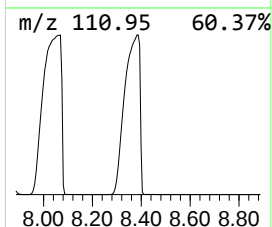
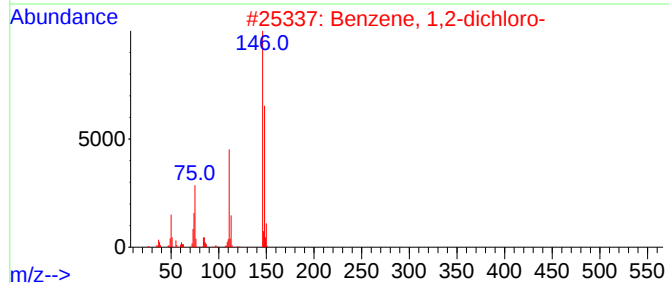
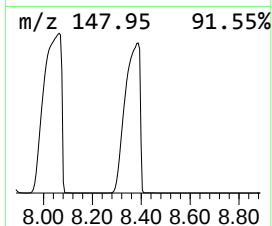
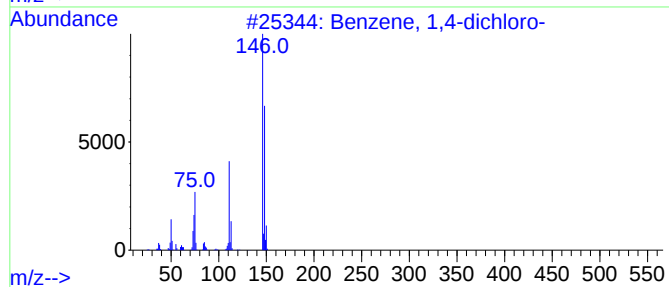
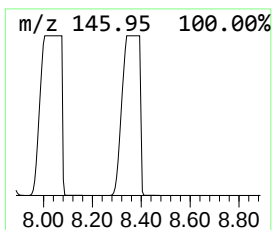
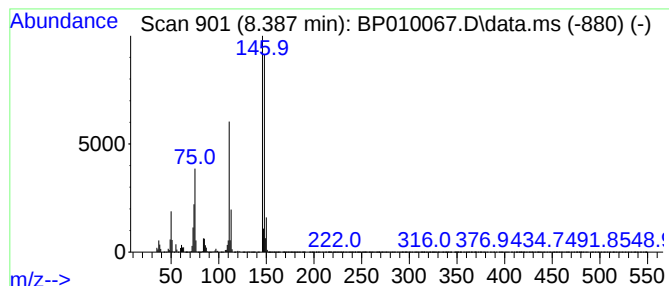
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Quant Title : SVOA CALIBRATION

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.
8.387	16.13 ng/ul	190795000	1,4-Dichlorobenzene-d4	7.993

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	98
2			Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	98
3			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	96
4			Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	95
5			Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	95



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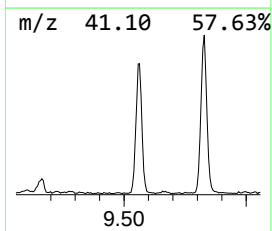
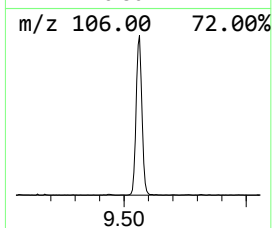
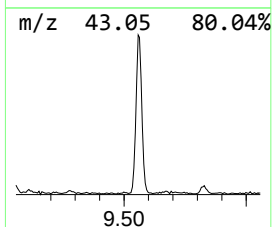
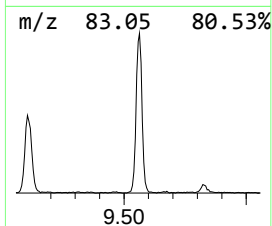
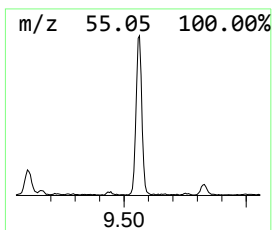
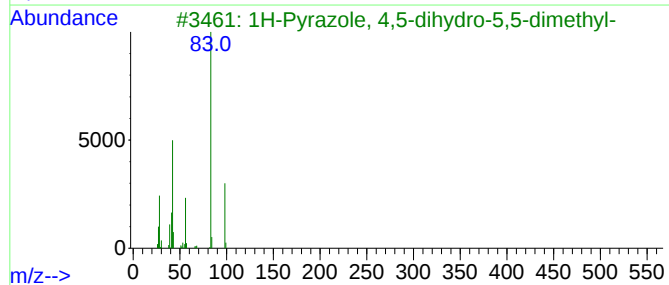
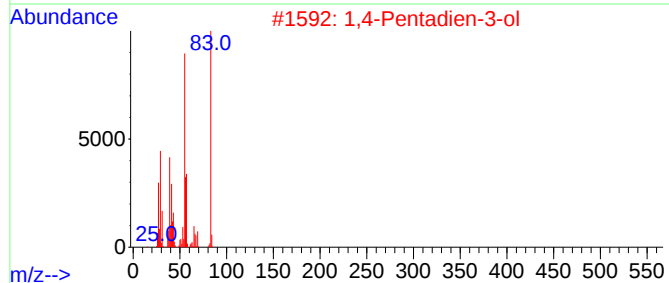
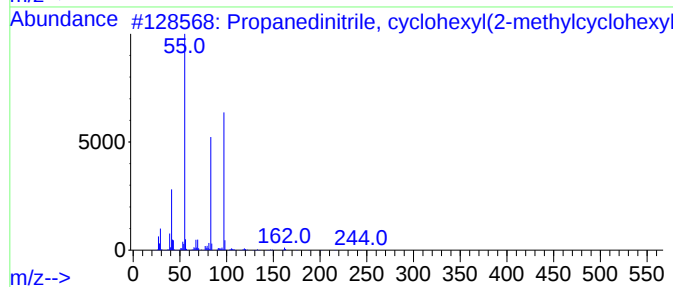
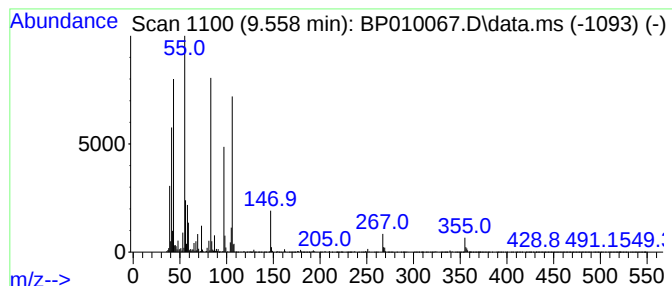
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

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

Peak Number 5 unknown-01 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.558	7.42 ng/ul	526234	Naphthalene-d8	10.758

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propanedinitrile, cyclohexyl(2-m...	244	C16H24N2	074764-55-9	38
2			1,4-Pentadien-3-ol	84	C5H8O	000922-65-6	30
3			1H-Pyrazole, 4,5-dihydro-5,5-dim...	98	C5H10N2	004320-85-8	27
4			Cyclopropane, 1,1,2,2-tetramethyl-	98	C7H14	004127-47-3	22
5			Aziridine, 1-(1-propenyl)-, (E)-	83	C5H9N	080839-91-4	22



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010067.D
Acq On : 22 Apr 2022 22:14
Operator : CG/JU
Sample : N2490-07
Misc :
ALS Vial : 63 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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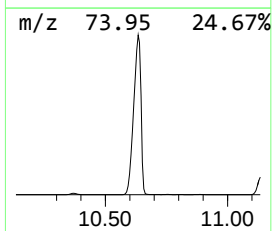
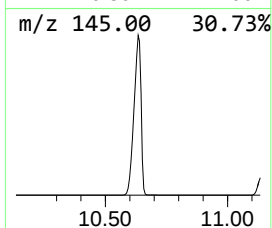
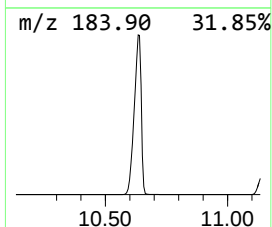
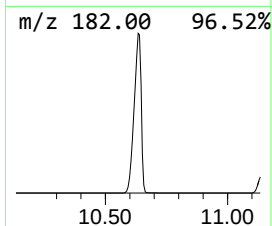
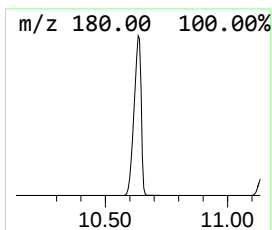
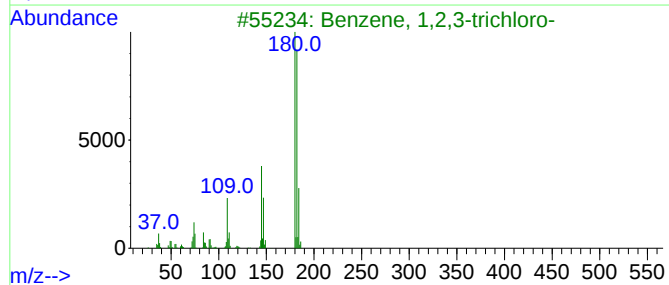
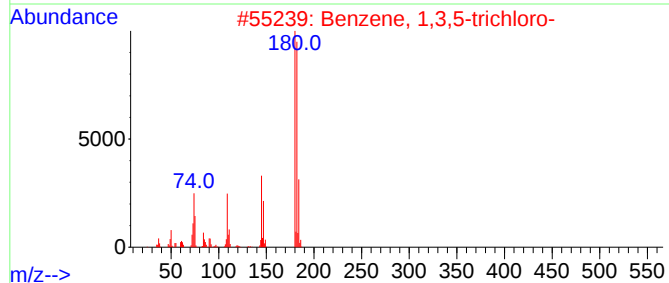
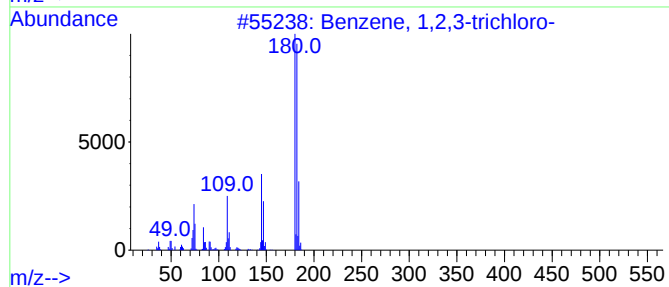
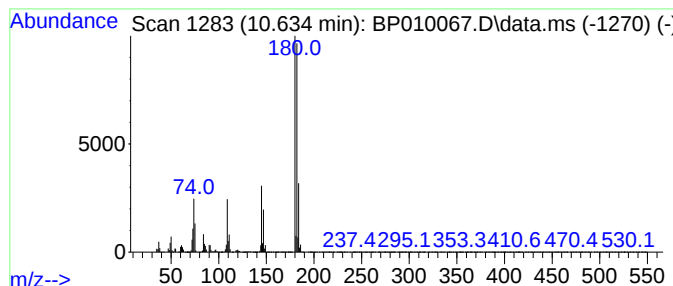
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Benzene, 1,2,3-trichloro- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.634	751.78 ng/ul	53308000	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, 1,2,4-trichloro-	180	C6H3Cl3	000120-82-1	97
5			Benzene, 1,3,5-trichloro-	180	C6H3Cl3	000108-70-3	97



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010067.D
Acq On : 22 Apr 2022 22:14
Operator : CG/JU
Sample : N2490-07
Misc :
ALS Vial : 63 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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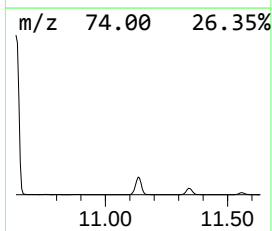
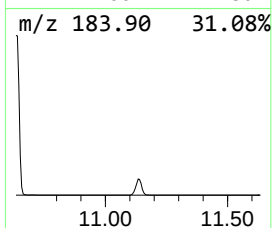
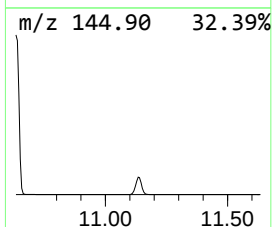
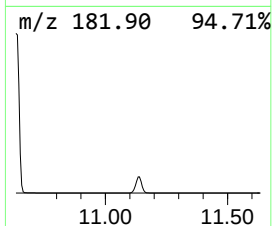
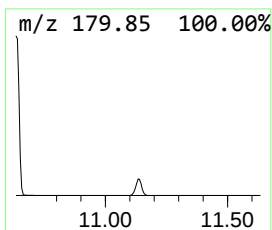
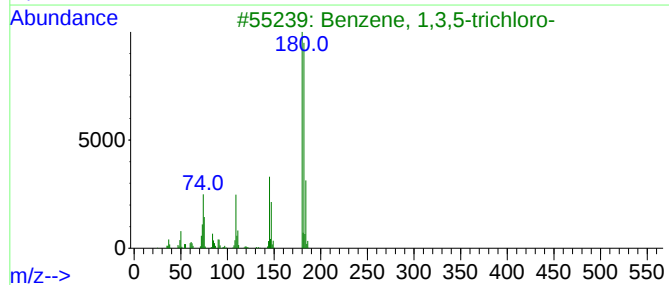
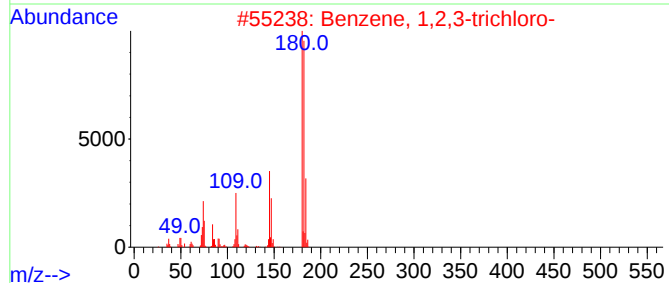
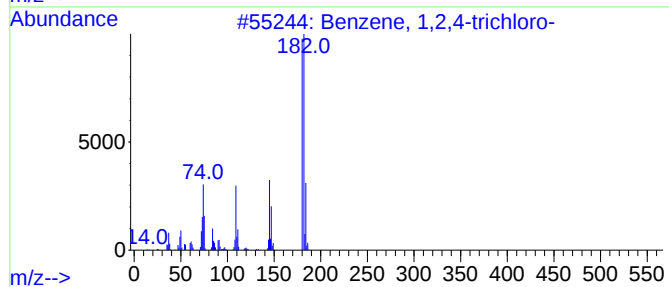
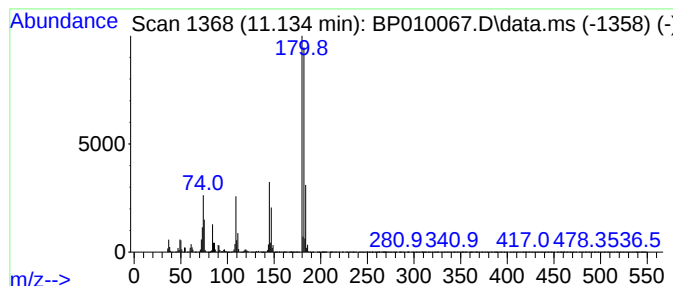
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Benzene, 1,2,4-trichloro- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.134	66.64 ng/ul	4725290	Naphthalene-d8	10.758

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,4-trichloro-	180	C6H3Cl3	000120-82-1	98
2			Benzene, 1,2,3-trichloro-	180	C6H3Cl3	000087-61-6	98
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 : BP010067.D
Acq On : 22 Apr 2022 22:14
Operator : CG/JU
Sample : N2490-07
Misc :
ALS Vial : 63 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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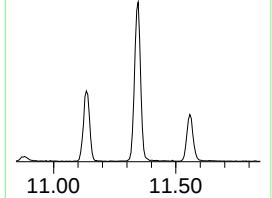
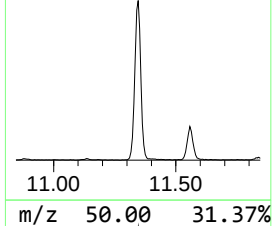
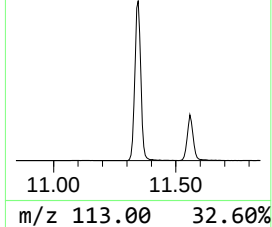
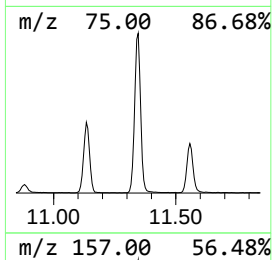
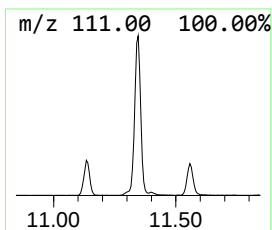
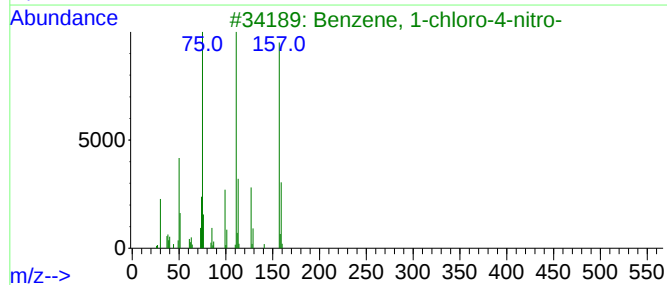
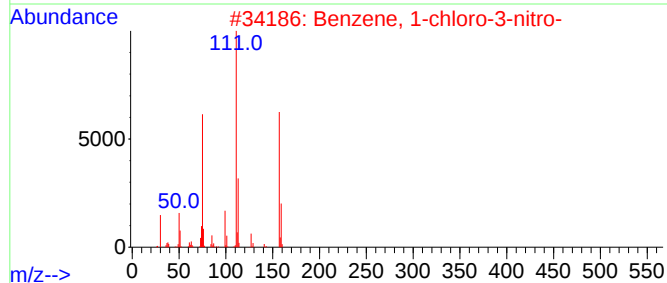
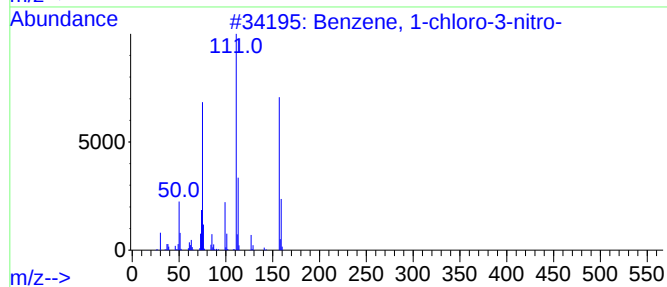
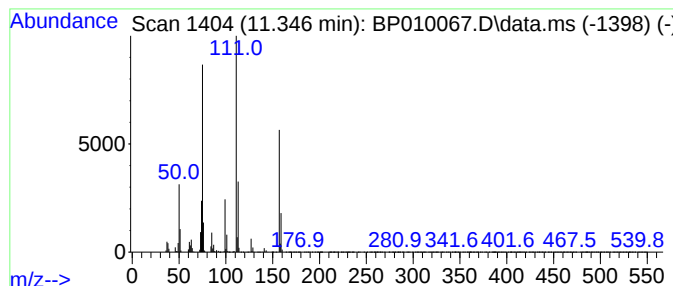
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Benzene, 1-chloro-3-nitro- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.346	27.29 ng/ul	1935040	Naphthalene-d8	10.758

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	95
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 : BP010067.D
Acq On : 22 Apr 2022 22:14
Operator : CG/JU
Sample : N2490-07
Misc :
ALS Vial : 63 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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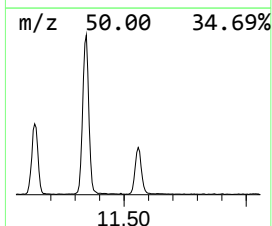
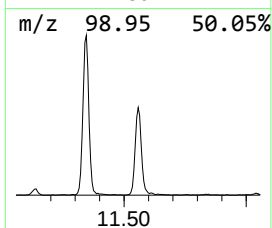
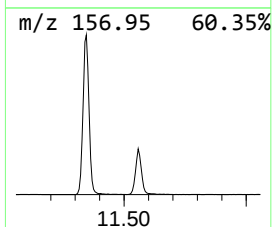
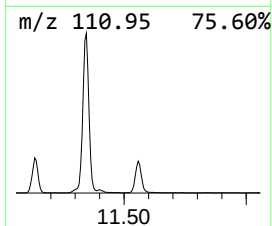
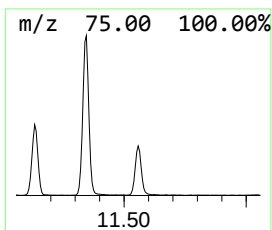
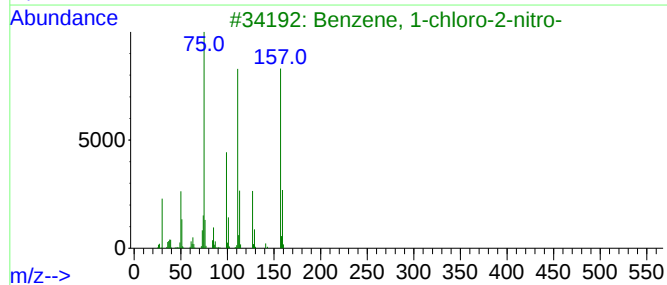
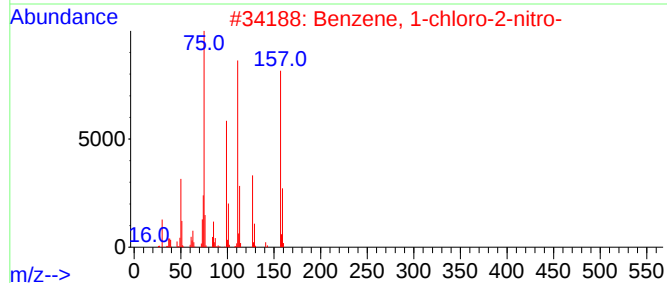
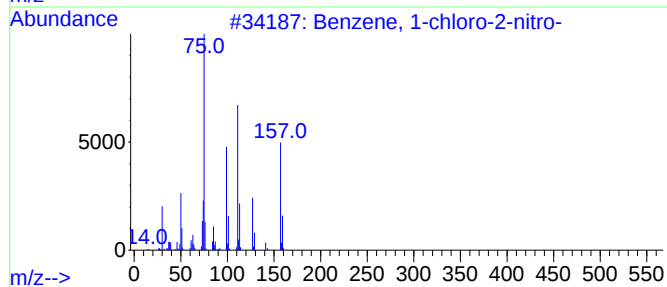
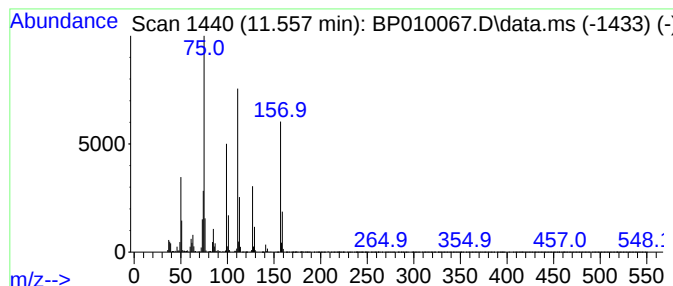
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Benzene, 1-chloro-2-nitro- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.557	8.70 ng/ul	616887	Naphthalene-d8	10.758

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-2-nitro-	157	C6H4ClNO2	000088-73-3	97
3			Benzene, 1-chloro-2-nitro-	157	C6H4ClNO2	000088-73-3	96
4			Benzene, 1-chloro-2-nitro-	157	C6H4ClNO2	000088-73-3	96
5			Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010067.D
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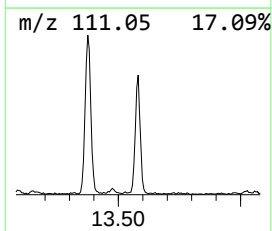
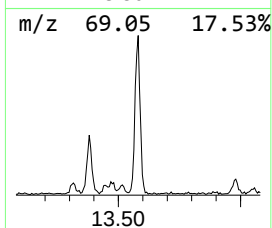
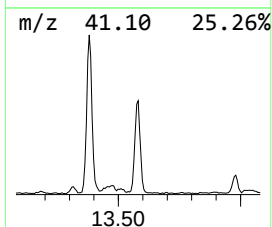
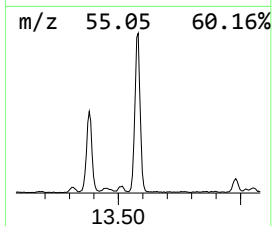
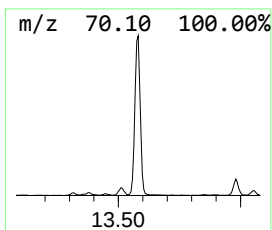
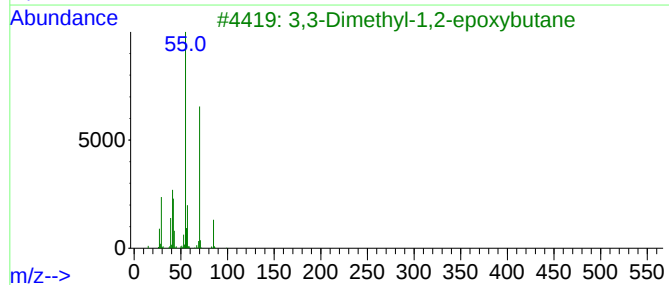
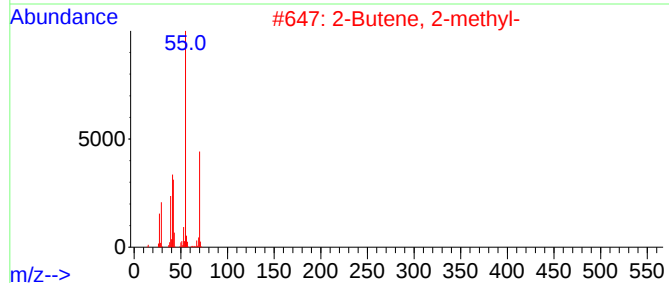
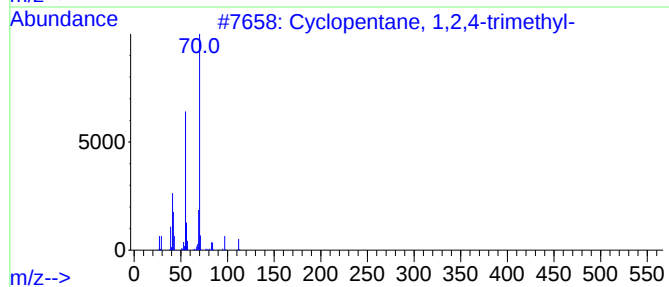
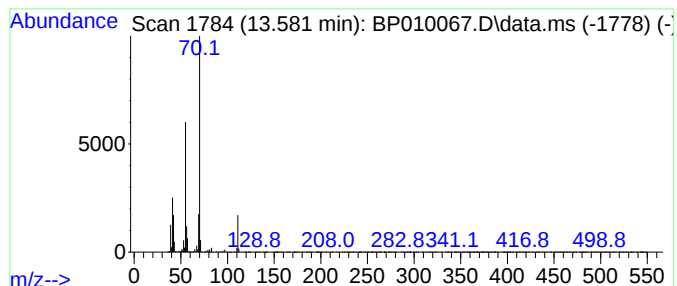
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 (DEL) Alkane: Cyclic13.581 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.581	4.51 ng/ul	493183	Acenaphthene-d10	14.569

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentane, 1,2,4-trimethyl-	112	C8H16	002815-58-9	72
2			2-Butene, 2-methyl-	70	C5H10	000513-35-9	72
3			3,3-Dimethyl-1,2-epoxybutane	100	C6H12O	002245-30-9	59
4			Cyclopentane, 1,2,4-trimethyl-	112	C8H16	002815-58-9	56
5			Cyclopentane, 1,2,4-trimethyl-, ...	112	C8H16	004850-28-6	56



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
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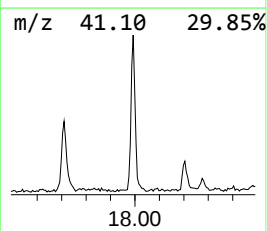
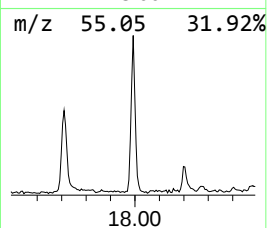
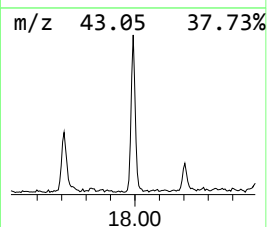
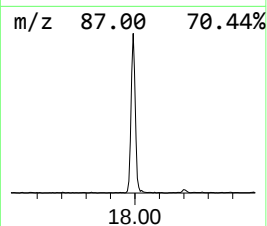
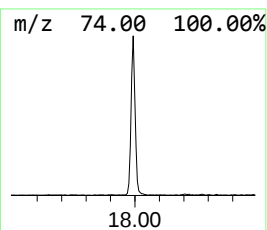
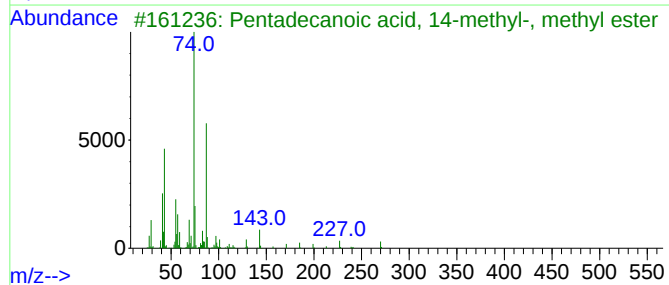
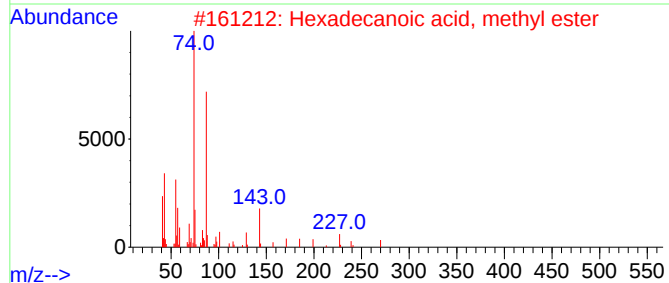
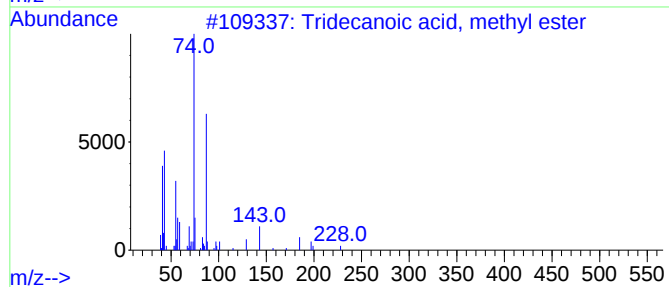
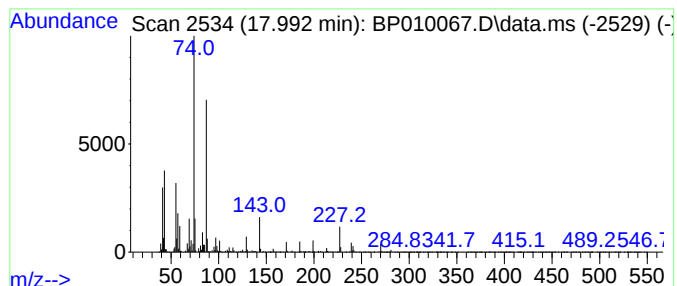
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 Tridecanoic acid, methyl ester Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.992	3.33 ng/ul	432848	Phenanthrene-d10	17.322

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	98
2			Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	98
3			Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	93
4			Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	91
5			Methyl tetradecanoate	242	C15H30O2	000124-10-7	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
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ALS Vial : 63 Sample Multiplier: 1

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ClientSampleId :
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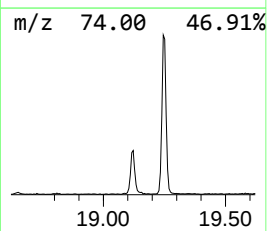
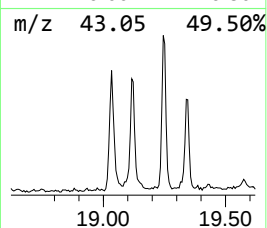
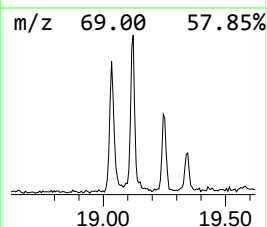
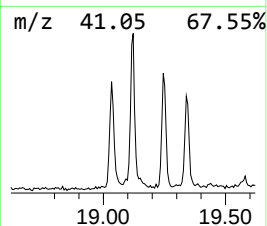
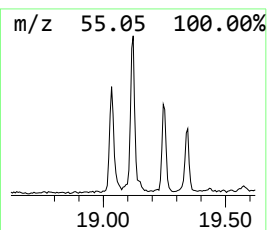
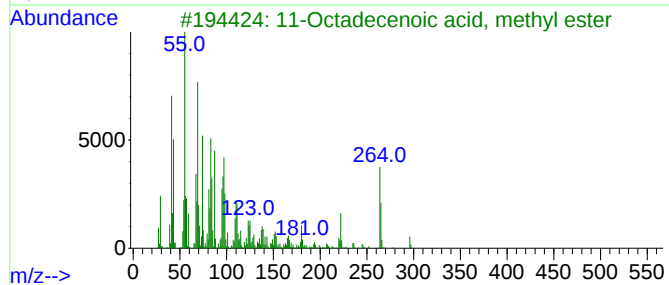
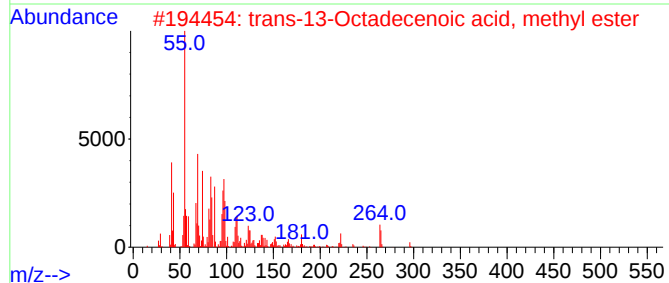
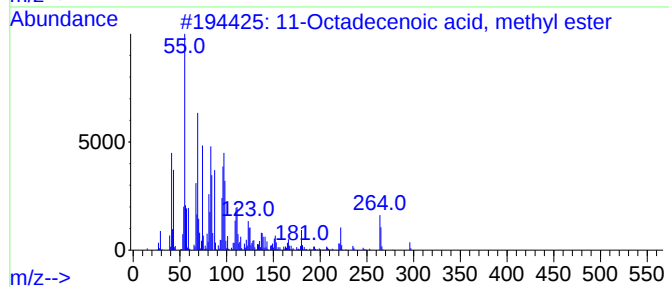
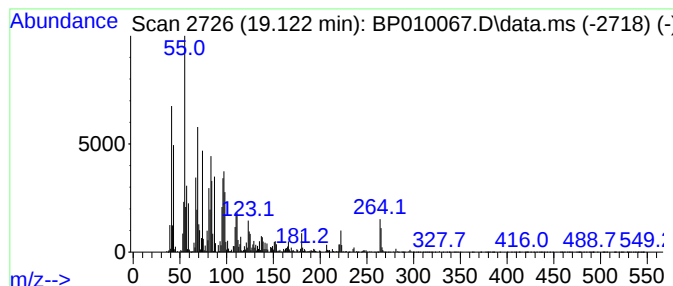
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 11-Octadecenoic acid, methyl... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.122	3.97 ng/ul	517234	Phenanthrene-d10	17.322

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			11-Octadecenoic acid, methyl ester	296	C19H36O2	052380-33-3	94
2			trans-13-Octadecenoic acid, methyl ester	296	C19H36O2	1000333-61-3	94
3			11-Octadecenoic acid, methyl ester	296	C19H36O2	052380-33-3	94
4			9-Octadecenoic acid, methyl ester	296	C19H36O2	001937-62-8	94
5			cis-13-Octadecenoic acid, methyl ester	296	C19H36O2	1010333-58-3	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010067.D
Acq On : 22 Apr 2022 22:14
Operator : CG/JU
Sample : N2490-07
Misc :
ALS Vial : 63 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0J44

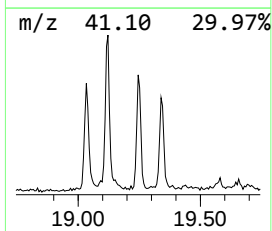
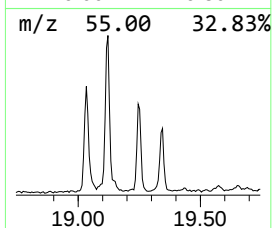
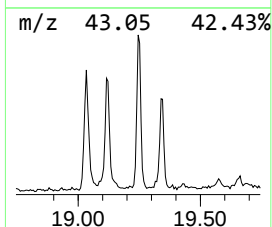
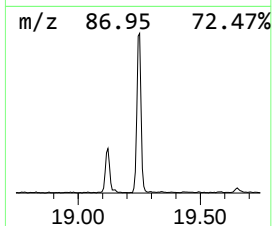
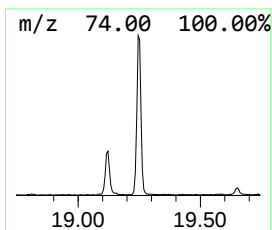
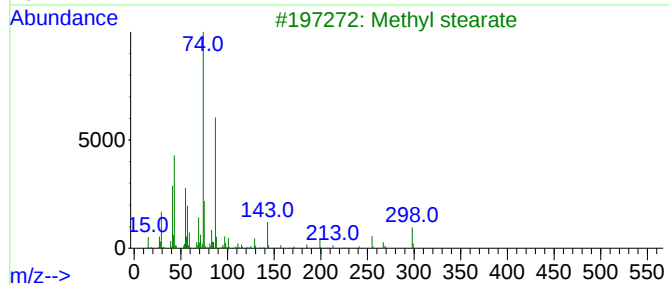
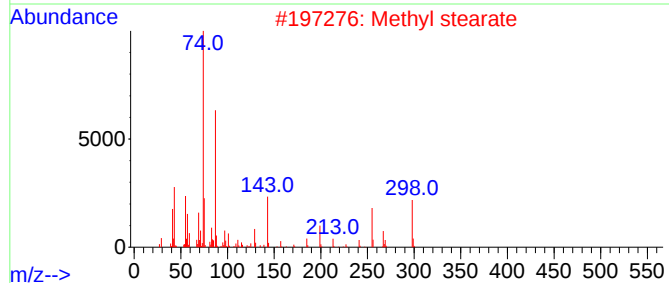
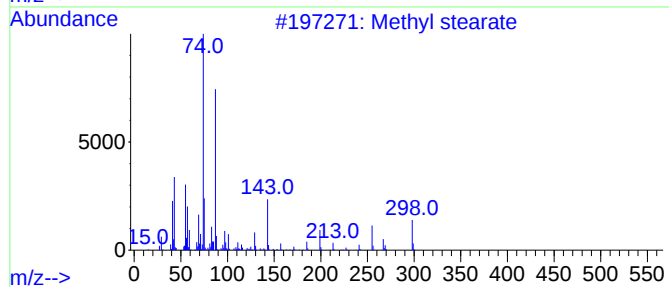
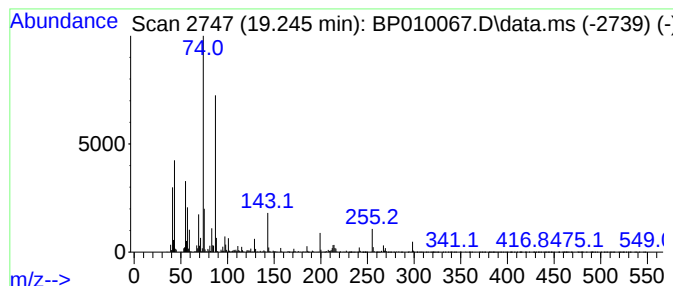
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 Methyl stearate Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.245	3.51 ng/ul	456454	Phenanthrene-d10	17.322

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methyl stearate	298	C19H38O2	000112-61-8	98
2			Methyl stearate	298	C19H38O2	000112-61-8	97
3			Methyl stearate	298	C19H38O2	000112-61-8	95
4			Methyl stearate	298	C19H38O2	000112-61-8	90
5			Methyl stearate	298	C19H38O2	000112-61-8	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
 Data File : BP010067.D
 Acq On : 22 Apr 2022 22:14
 Operator : CG/JU
 Sample : N2490-07
 Misc :
 ALS Vial : 63 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0J44

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-	5.364	29.1	ng/ul	343991000	1	7.993	236568000	20.0
Benzene, 1,2-di...	7.852	2.1	ng/ul	24516000	1	7.993	236568000	20.0
Benzene, 1,4-di...	8.063	20.0	ng/ul	236568000	1	7.993	236568000	20.0
Benzene, 1,3-di...	8.387	16.1	ng/ul	190795000	1	7.993	236568000	20.0
unknown-01	9.558	7.4	ng/ul	526234	2	10.758	1418180	20.0
Benzene, 1,2,3-...	10.634	751.8	ng/ul	53308000	2	10.758	1418180	20.0
Benzene, 1,2,4-...	11.134	66.6	ng/ul	4725290	2	10.758	1418180	20.0
Benzene, 1-chlo...	11.346	27.3	ng/ul	1935040	2	10.758	1418180	20.0
Benzene, 1-chlo...	11.557	8.7	ng/ul	616887	2	10.758	1418180	20.0
(DEL) Alkane: C...	13.581	4.5	ng/ul	493183	3	14.569	2184740	20.0
Tridecanoic aci...	17.992	3.3	ng/ul	432848	4	17.322	2602700	20.0
11-Octadecenoic...	19.122	4.0	ng/ul	517234	4	17.322	2602700	20.0
Methyl stearate	19.245	3.5	ng/ul	456454	4	17.322	2602700	20.0