

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM011422\
 Data File : BM033736.D
 Acq On : 14 Jan 2022 20:00
 Operator : CG/JU
 Sample : N1121-03
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 WASTE-CLASS-LAUNDRY-PIT

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Filtering: 5

Sampling : 1

Min Area: 3 % 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_M\Methods\8270-BM010522.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BM033736.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.360	17	21	28	rVB	275706	297681	3.80%	0.617%
2	4.543	216	222	230	rBV	317393	413068	5.28%	0.856%
3	5.013	295	302	304	rBV	187588	259913	3.32%	0.539%
4	5.037	304	306	314	rVB2	178982	224716	2.87%	0.466%
5	5.513	380	387	395	rBV	3110631	4390274	56.08%	9.097%
6	7.125	653	661	668	rBV	2657951	4221229	53.92%	8.746%
7	7.966	795	804	810	rBV	615680	1002009	12.80%	2.076%
8	9.131	990	1002	1012	rBV	2052864	3508532	44.81%	7.270%
9	10.778	1272	1282	1288	rBV	855053	1439455	18.39%	2.983%
10	13.230	1692	1699	1711	rBV	5352910	7829103	100.00%	16.222%
11	14.601	1925	1932	1947	rBV	1272576	1874702	23.95%	3.884%
12	15.601	2096	2102	2107	rBV	112045	162129	2.07%	0.336%
13	16.083	2177	2184	2195	rVB	3637782	5289299	67.56%	10.959%
14	16.336	2221	2227	2232	rVB	118561	164215	2.10%	0.340%
15	16.424	2237	2242	2246	rBV	266540	353861	4.52%	0.733%
16	16.483	2246	2252	2258	rVB2	264049	508677	6.50%	1.054%
17	16.542	2258	2262	2266	rBV2	285497	380336	4.86%	0.788%
18	16.589	2266	2270	2275	rVB	182116	232213	2.97%	0.481%
19	16.665	2280	2283	2285	rVV	93116	102802	1.31%	0.213%
20	16.689	2285	2287	2291	rVB	80412	85941	1.10%	0.178%
21	16.742	2291	2296	2303	rVB3	255335	464478	5.93%	0.962%
22	16.807	2303	2307	2311	rBV	269208	375302	4.79%	0.778%
23	16.883	2317	2320	2327	rVB	156814	208646	2.67%	0.432%
24	17.336	2391	2397	2408	rBV	1277272	1820076	23.25%	3.771%
25	18.230	2544	2549	2553	rBV	112199	168778	2.16%	0.350%
26	19.430	2749	2753	2761	rVB5	72716	119949	1.53%	0.249%
27	19.501	2762	2765	2773	rVB2	94648	122037	1.56%	0.253%
28	19.948	2835	2841	2847	rBV	6568245	7666467	97.92%	15.885%
29	20.124	2866	2871	2872	rBV4	100699	143791	1.84%	0.298%
30	20.224	2883	2888	2893	rBV2	263014	509443	6.51%	1.056%
31	21.494	3100	3104	3109	rBV	1263201	1571029	20.07%	3.255%
32	22.071	3199	3202	3209	rVB2	327596	553331	7.07%	1.146%
33	23.912	3509	3515	3526	rVB2	881546	1799615	22.99%	3.729%

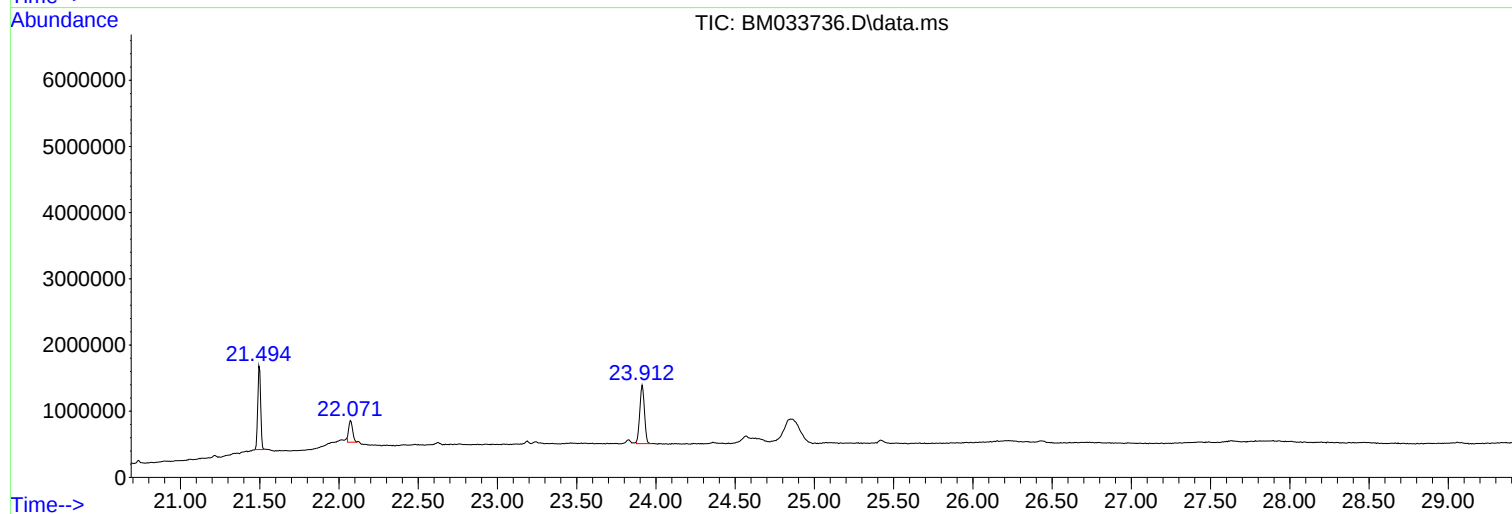
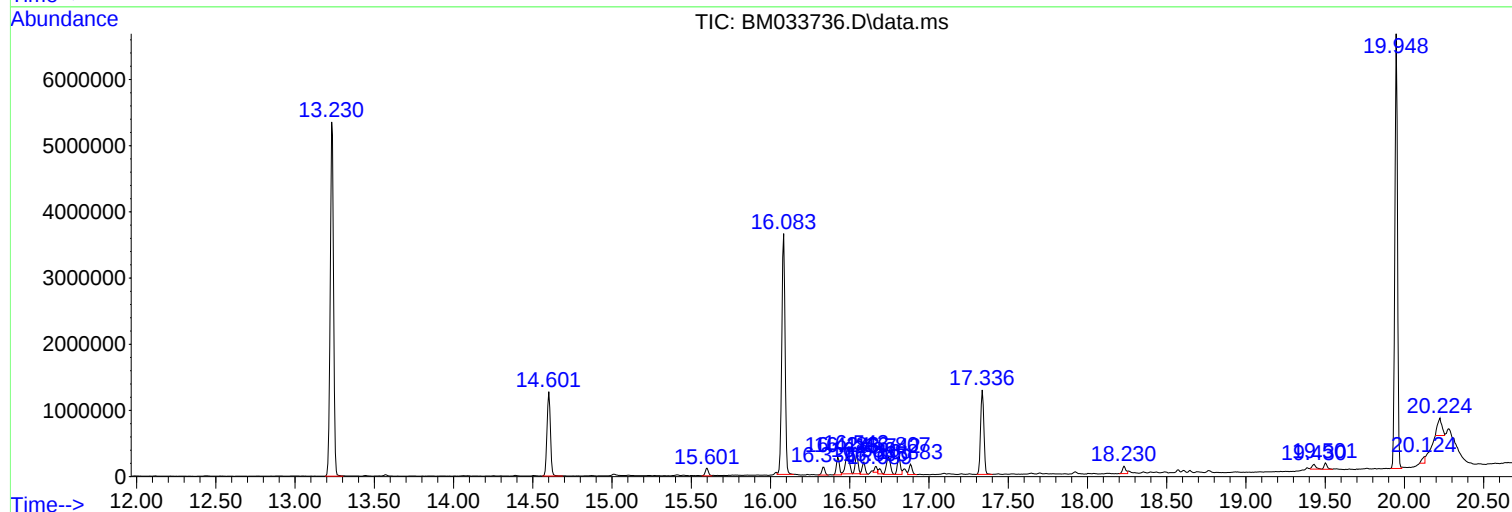
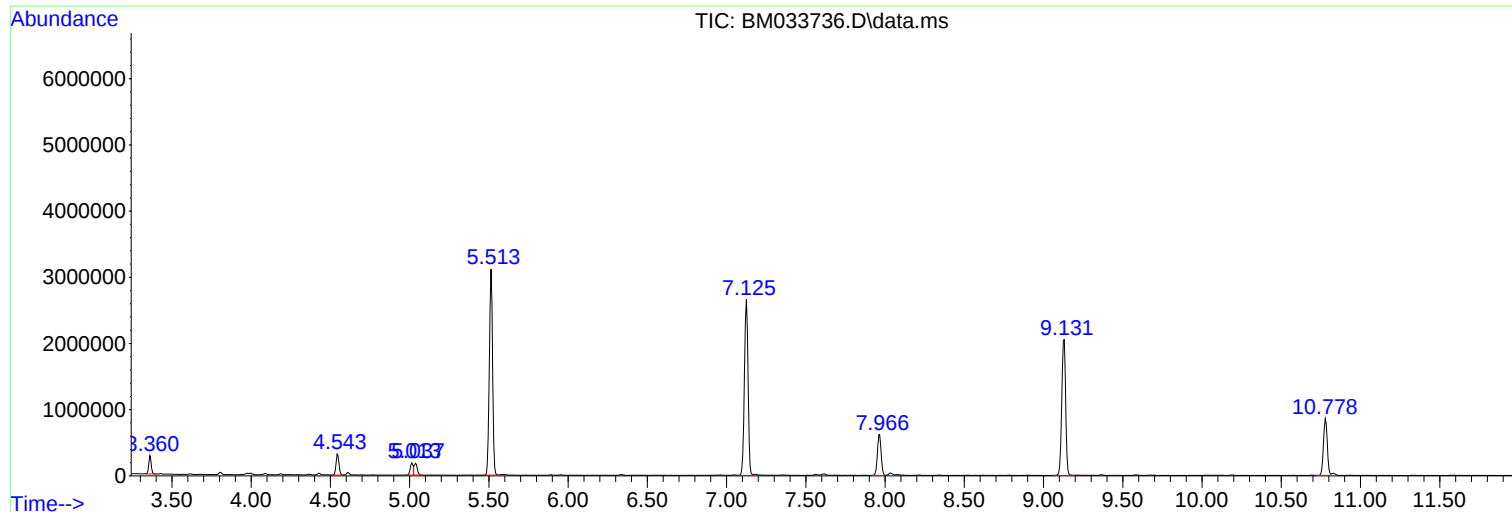
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM010522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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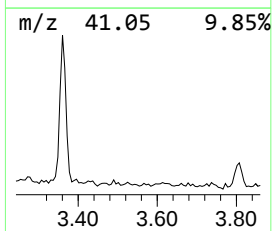
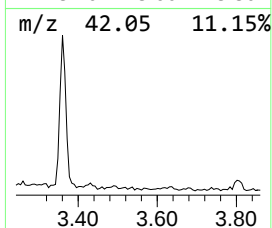
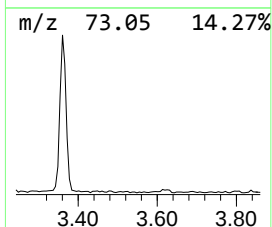
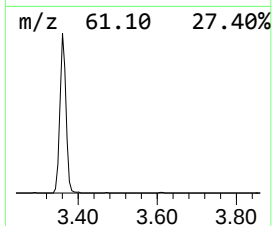
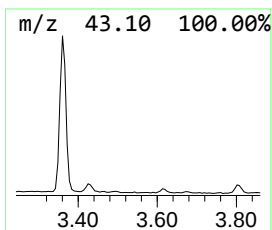
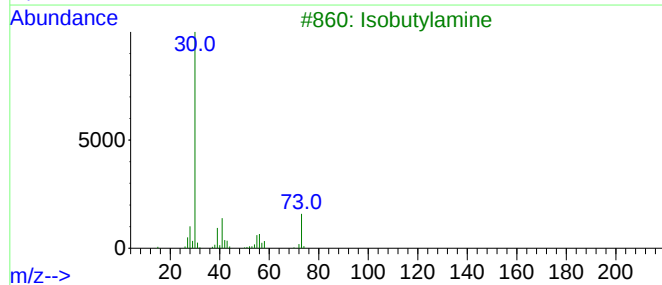
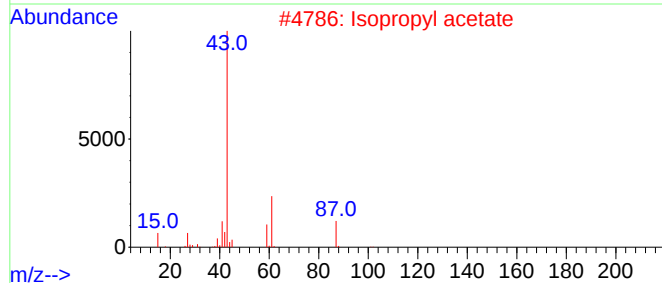
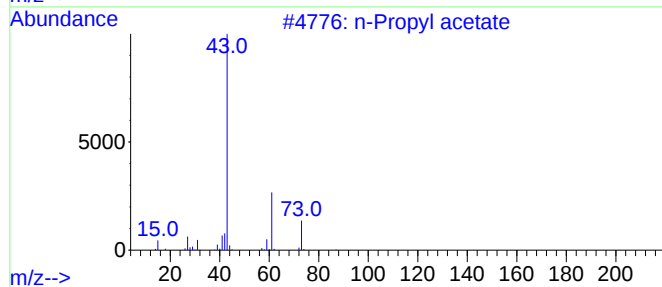
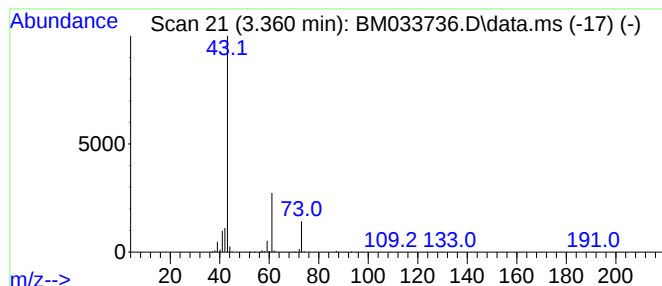
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM010522.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 n-Propyl acetate Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.360	5.94 ng	297681	1,4-Dichlorobenzene-d4	7.966

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Propyl acetate	102	C5H10O2	000109-60-4	83
2			Isopropyl acetate	102	C5H10O2	000108-21-4	38
3			Isobutylamine	73	C4H11N	000078-81-9	9
4			Isobutyl acetate	116	C6H12O2	000110-19-0	9
5			1-Butanamine	73	C4H11N	000109-73-9	7



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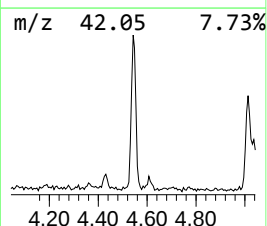
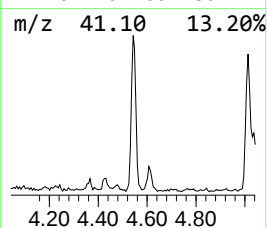
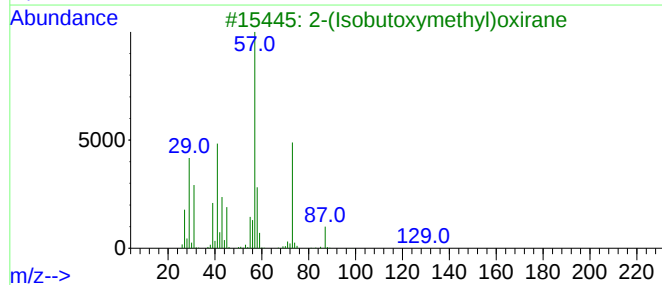
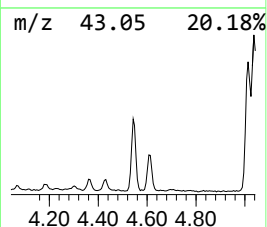
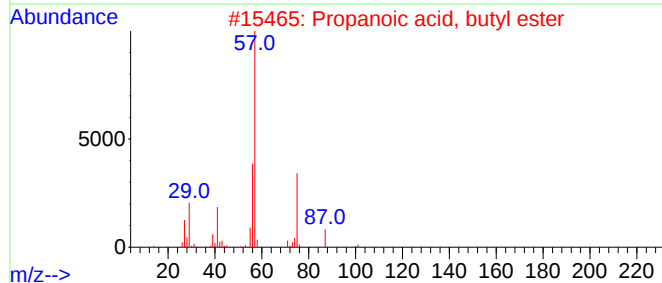
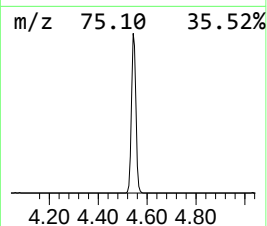
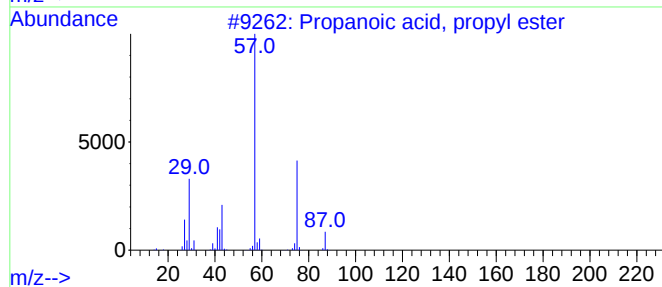
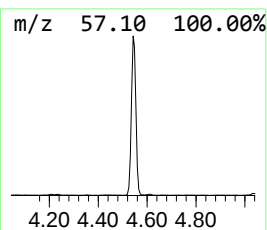
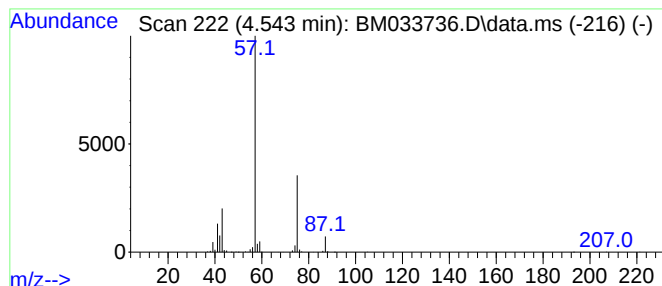
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM010522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
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Peak Number 2 Propanoic acid, propyl ester Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.543	8.24 ng	413068	1,4-Dichlorobenzene-d4	7.966

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propanoic acid, propyl ester	116	C6H12O2	000106-36-5	83
2			Propanoic acid, butyl ester	130	C7H14O2	000590-01-2	50
3			2-(Isobutoxymethyl)oxirane	130	C7H14O2	003814-55-9	9
4			di-tert-Butyl dicarbonate	218	C10H18O5	024424-99-5	9
5			1-Butanamine, 2-methyl-	87	C5H13N	000096-15-1	7



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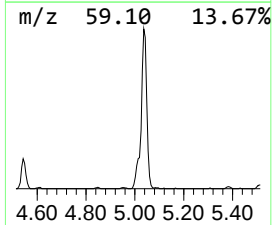
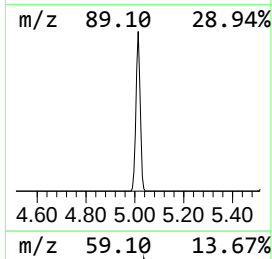
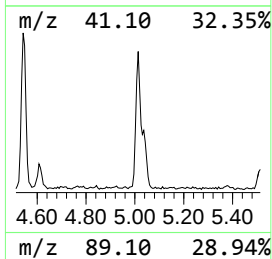
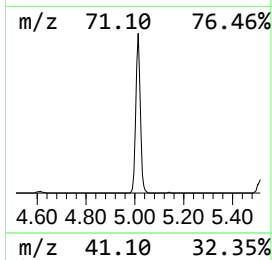
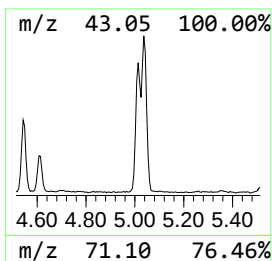
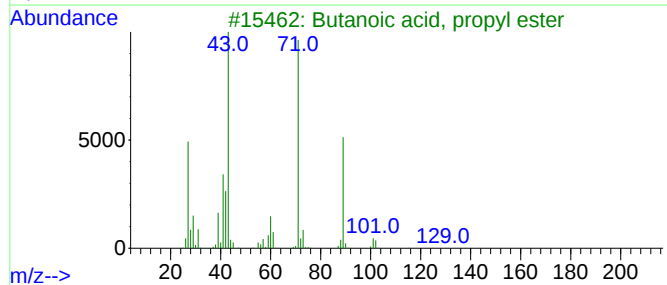
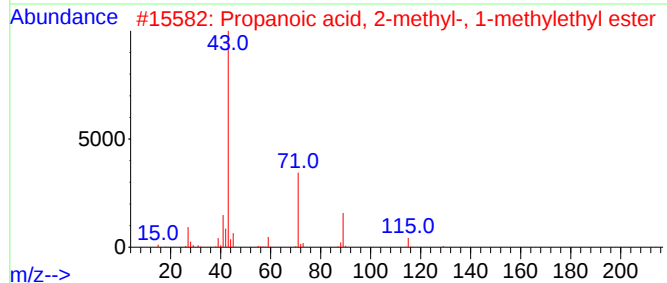
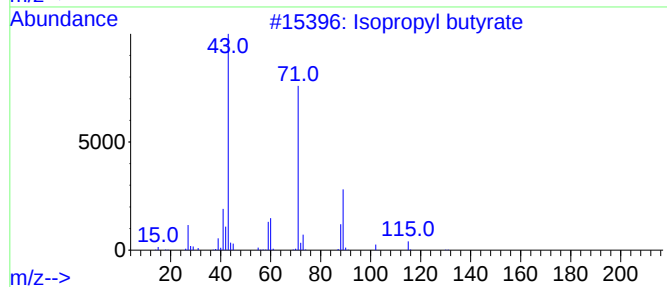
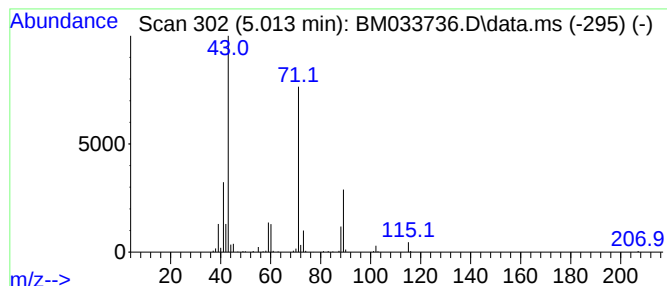
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Isopropyl butyrate Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.013	5.19 ng	259913	1,4-Dichlorobenzene-d4	7.966

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Isopropyl butyrate	130	C7H14O2	000638-11-9	90
2			Propanoic acid, 2-methyl-, 1-met...	130	C7H14O2	000617-50-5	72
3			Butanoic acid, propyl ester	130	C7H14O2	000105-66-8	64
4			Ethylene glycol Formate Isobutyrate	160	C7H12O4	1000458-48-0	64
5			Propanoic acid, 2-methyl-, propy...	130	C7H14O2	000644-49-5	50



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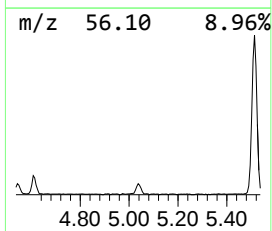
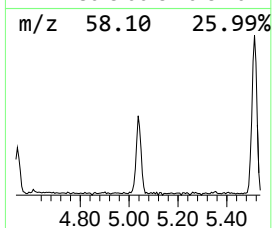
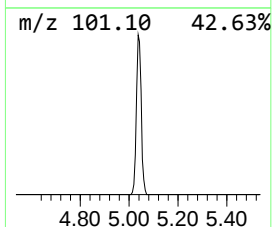
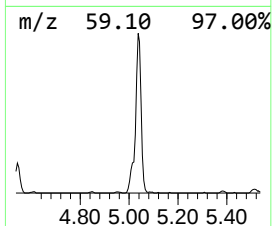
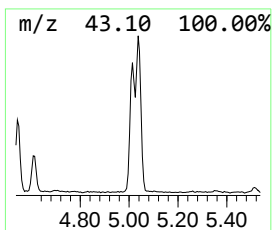
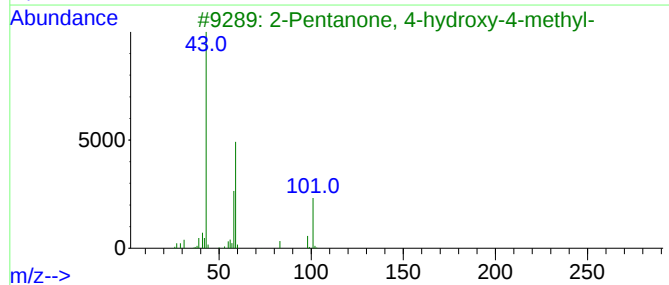
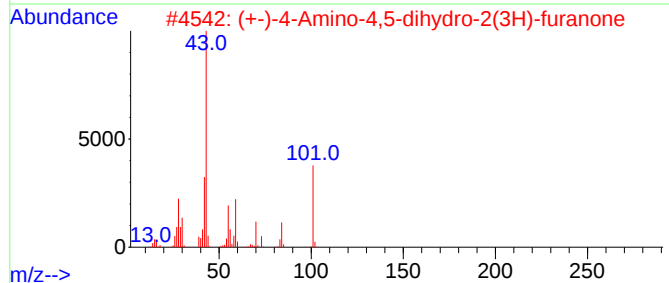
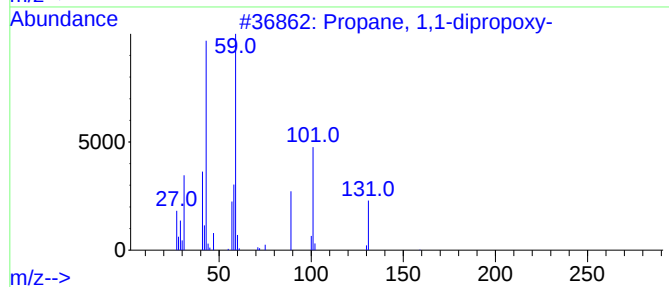
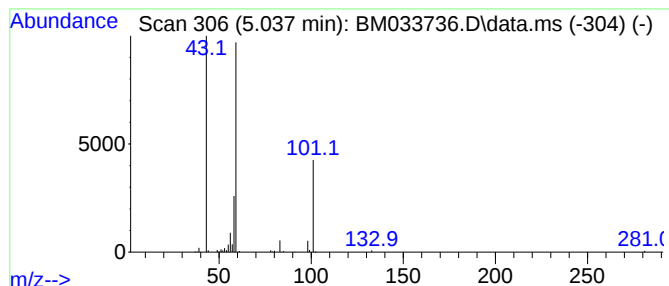
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 unknown5.037 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.037	4.49 ng	224716	1,4-Dichlorobenzene-d4	7.966

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propane, 1,1-dipropoxy-	160	C9H20O2	004744-11-0	40
2			(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	9
3			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	9
4			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	5
5			Propanamide, N-ethyl-	101	C5H11NO	005129-72-6	5



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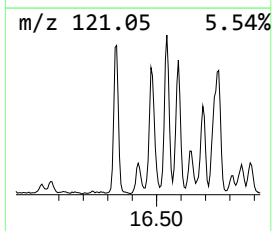
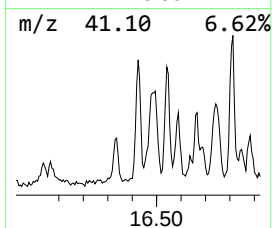
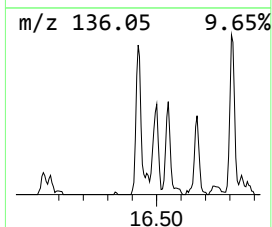
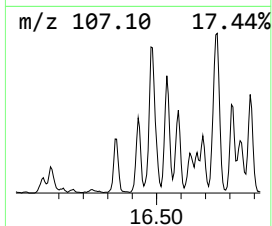
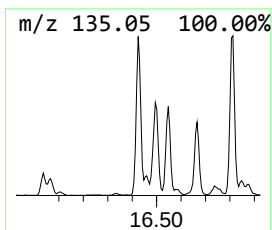
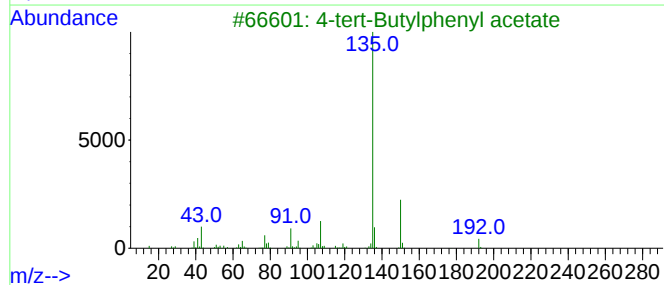
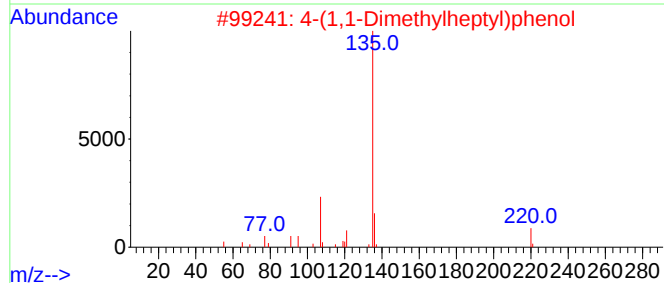
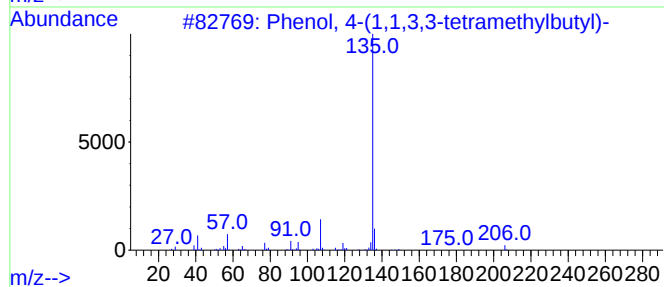
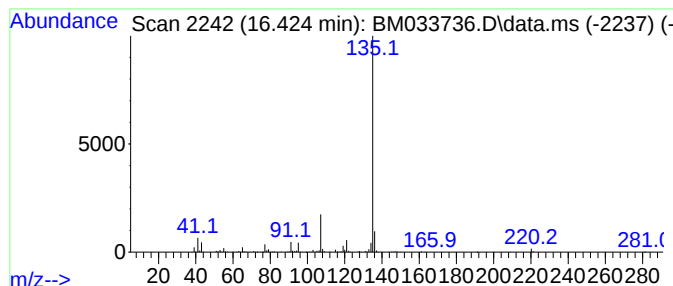
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Phenol, 4-(1,1,3,3-tetramet... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.424	3.89 ng	353861	Phenanthrene-d10	17.336

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 4-(1,1,3,3-tetramethylbu...	206	C14H22O	000140-66-9	86
2			4-(1,1-Dimethylheptyl)phenol	220	C15H24O	1000384-80-3	83
3			4-tert-Butylphenyl acetate	192	C12H16O2	003056-64-2	78
4			Phenol, 2-(1,1,3,3-tetramethylbu...	206	C14H22O	003884-95-5	78
5			2-tert-Butylphenol, O-(n-propylo...	236	C14H20O3	1000487-37-9	78



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM011422\
Data File : BM033736.D
Acq On : 14 Jan 2022 20:00
Operator : CG/JU
Sample : N1121-03
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
WASTE-CLASS-LAUNDRY-PIT

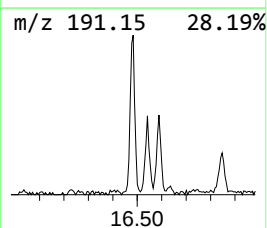
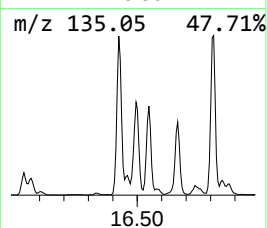
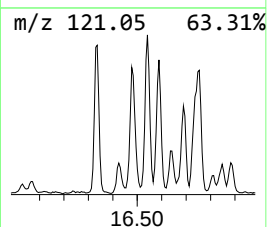
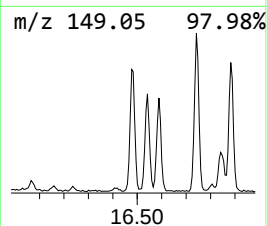
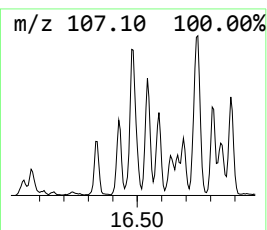
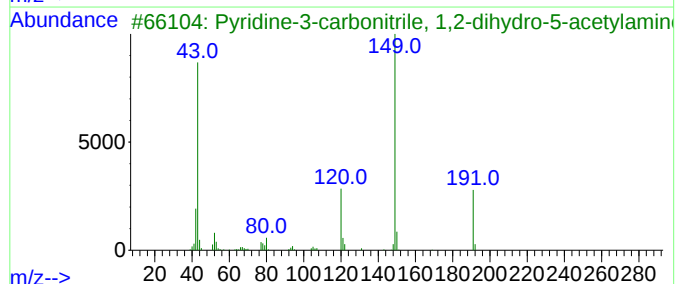
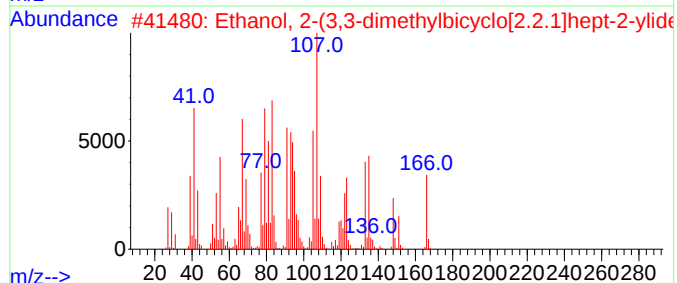
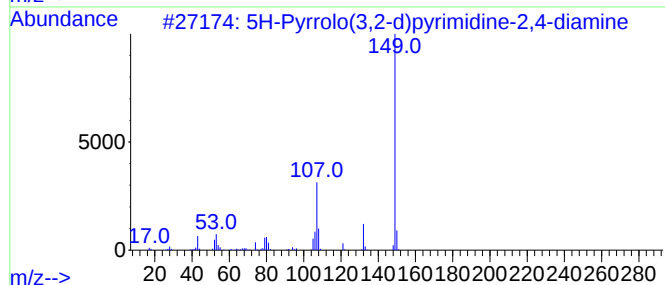
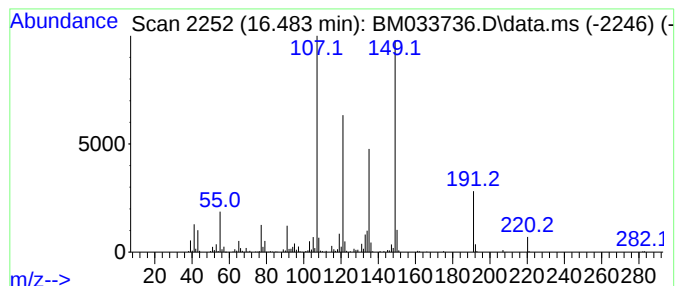
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 5H-Pyrrolo(3,2-d)pyrimidine... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.483	5.59 ng	508677	Phenanthrene-d10	17.336

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			5H-Pyrrolo(3,2-d)pyrimidine-2,4-...	149	C6H7N5	1000244-21-4	50
2			Ethanol, 2-(3,3-dimethylbicyclo[...]	166	C11H18O	002226-05-3	38
3			Pyridine-3-carbonitrile, 1,2-dih...	191	C9H9N3O2	220098-92-0	38
4			Phthalic acid, 4-chloro-3-methyl...	332	C18H17ClO4	1000309-84-2	35
5			Phthalic acid, 2-(methylthio)phe...	330	C18H18O4S	1000415-55-9	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM011422\
Data File : BM033736.D
Acq On : 14 Jan 2022 20:00
Operator : CG/JU
Sample : N1121-03
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
WASTE-CLASS-LAUNDRY-PIT

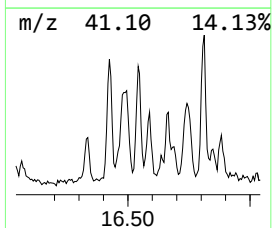
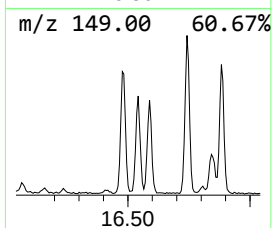
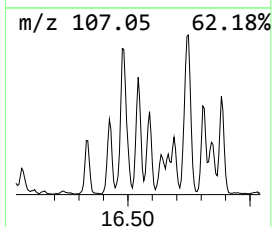
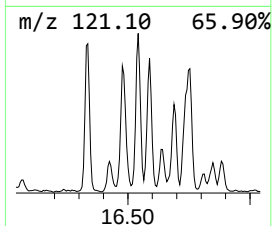
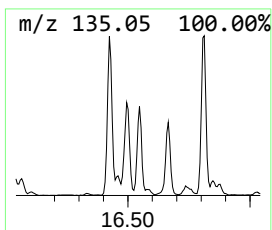
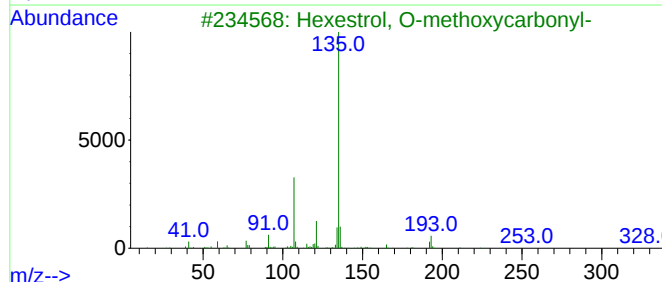
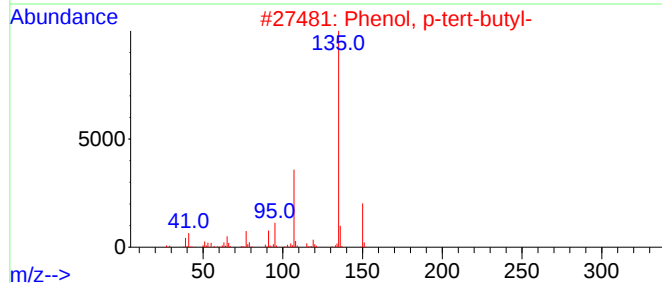
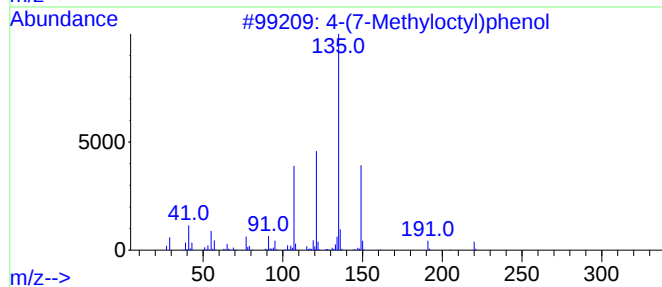
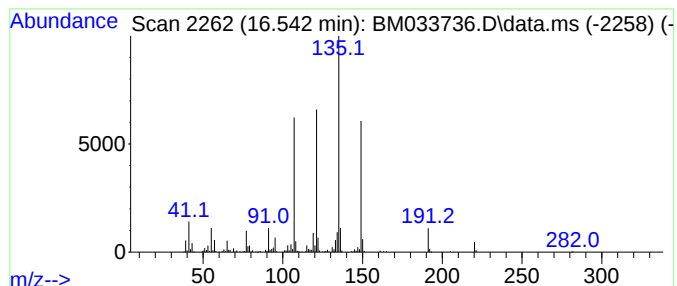
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 4-(7-Methyloctyl)phenol Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.542	4.18 ng	380336	Phenanthrene-d10	17.336

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		4-(7-Methyloctyl)phenol	220	C15H24O	024518-48-7	96
2		Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	64
3		Hexestrol, O-methoxycarbonyl-	328	C20H24O4	1000487-66-3	53
4		Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	53
5		Hexestrol, O-acetyl-	312	C20H24O3	1000374-87-8	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM011422\
Data File : BM033736.D
Acq On : 14 Jan 2022 20:00
Operator : CG/JU
Sample : N1121-03
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
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ClientSampleId :
WASTE-CLASS-LAUNDRY-PIT

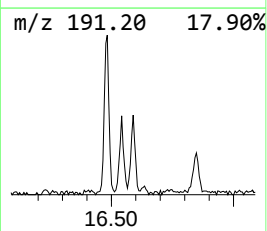
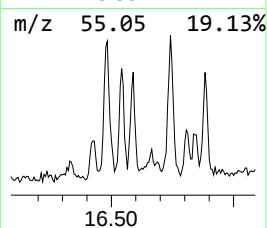
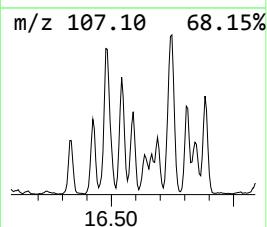
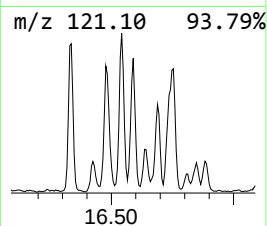
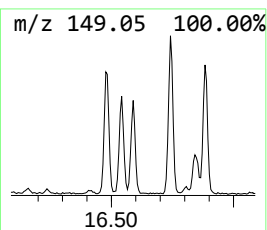
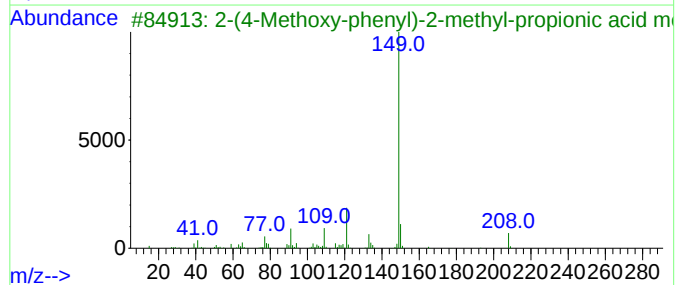
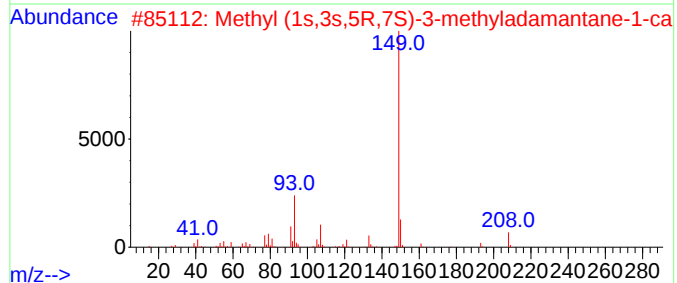
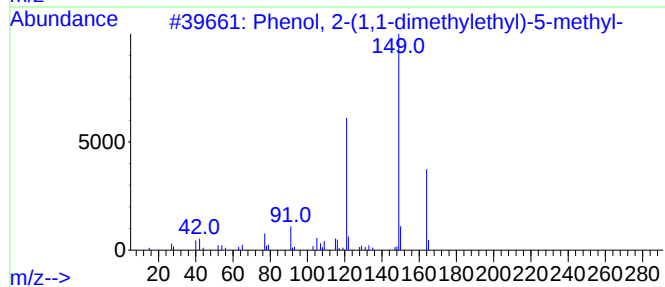
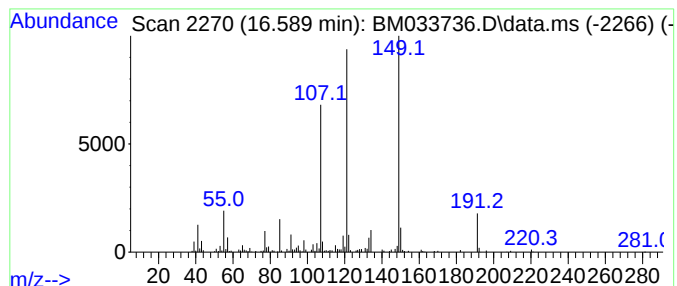
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Phenol, 2-(1,1-dimethylethy... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.589	2.55 ng	232213	Phenanthrene-d10	17.336

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 2-(1,1-dimethylethyl)-5-...	164	C11H16O	000088-60-8	50
2			Methyl (1s,3s,5R,7S)-3-methylada...	208	C13H20O2	1010504-39-1	35
3			2-(4-Methoxy-phenyl)-2-methyl-pr...	208	C12H16O3	006274-50-6	27
4			1-(2,6-Dimethyl-4-propoxyphenyl)...	220	C14H20O2	1000190-22-3	27
5			Phenol, 2-(1-methylpropyl)-	150	C10H14O	000089-72-5	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM011422\
Data File : BM033736.D
Acq On : 14 Jan 2022 20:00
Operator : CG/JU
Sample : N1121-03
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
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ClientSampleId :
WASTE-CLASS-LAUNDRY-PIT

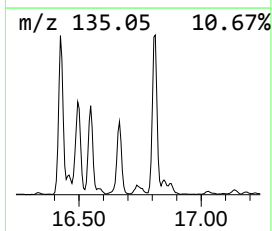
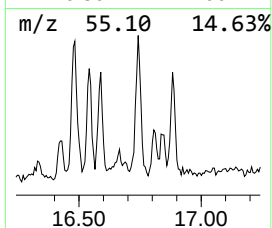
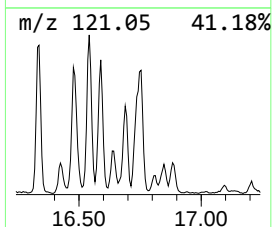
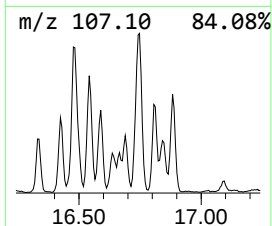
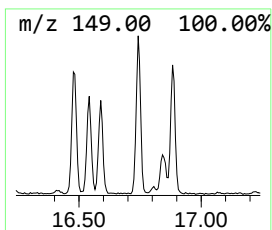
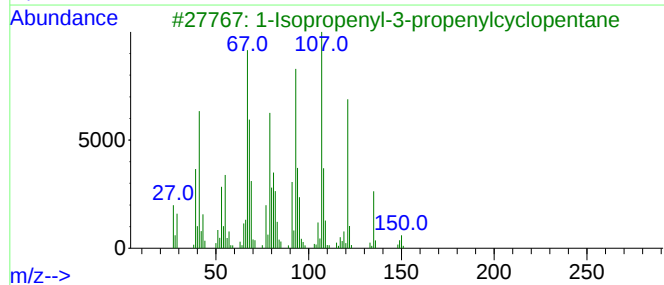
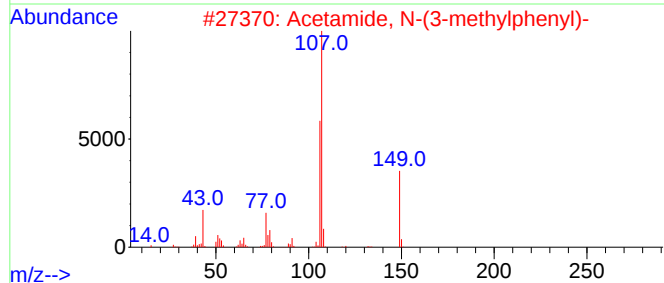
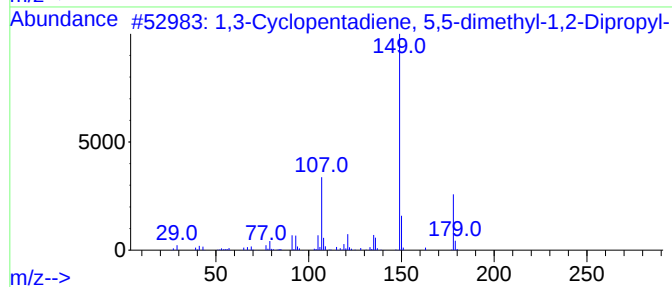
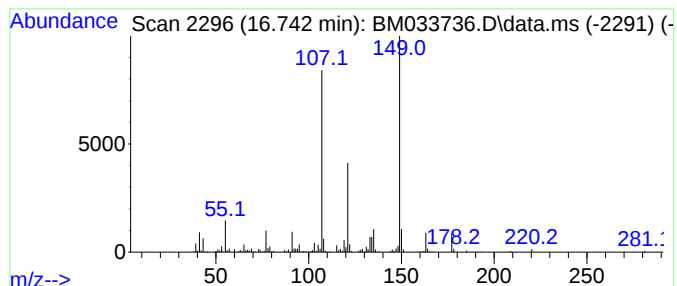
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 1,3-Cyclopentadiene, 5,5-di... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.742	5.10 ng	464478	Phenanthrene-d10	17.336

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1,3-Cyclopentadiene, 5,5-dimethy...	178	C13H22	1000163-88-0	72
2		Acetamide, N-(3-methylphenyl)-	149	C9H11NO	000537-92-8	47
3		1-Isopropenyl-3-propenylcyclopentane	150	C11H18	1000192-81-2	38
4		3,4-(methylenedioxy)-phenylbutan...	192	C11H12O3	1000378-99-2	38
5		2-tert-Butyl-6-methylphenol, n-b...	220	C15H24O	1000395-19-1	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM011422\
Data File : BM033736.D
Acq On : 14 Jan 2022 20:00
Operator : CG/JU
Sample : N1121-03
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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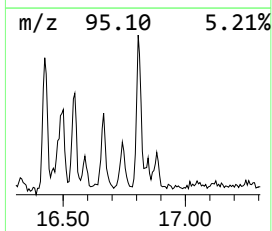
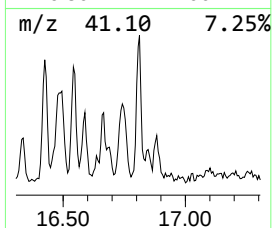
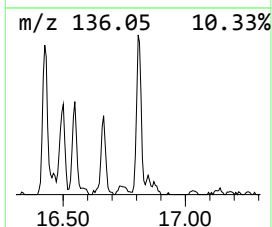
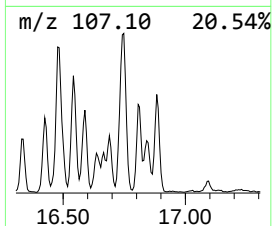
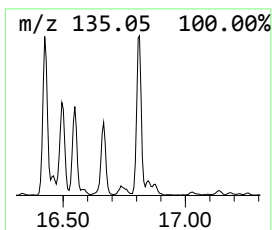
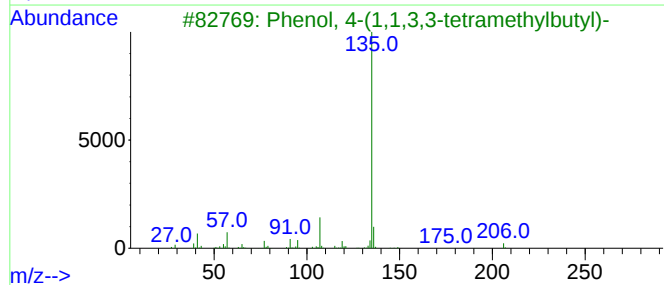
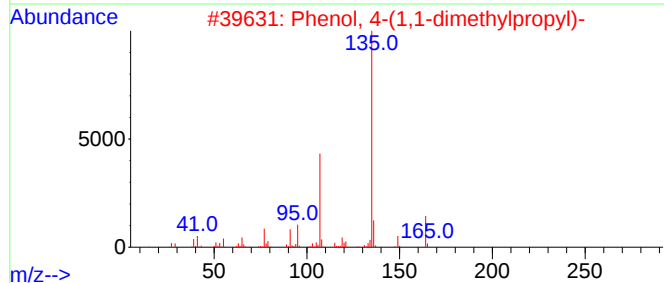
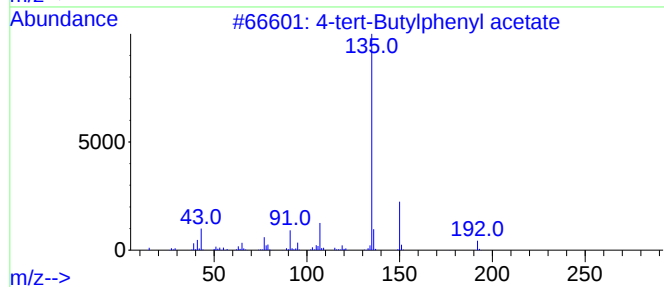
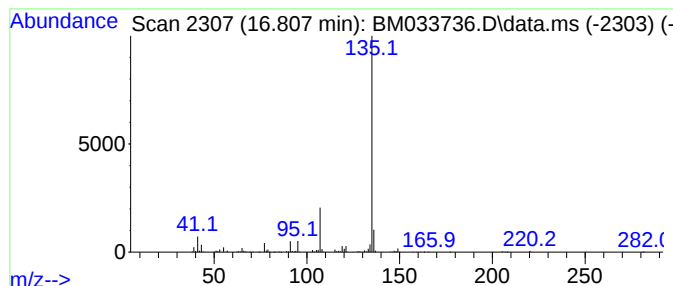
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 4-tert-Butylphenyl acetate Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.807	4.12 ng	375302	Phenanthrene-d10	17.336

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-tert-Butylphenyl acetate	192	C12H16O2	003056-64-2	78
2			Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	78
3			Phenol, 4-(1,1,3,3-tetramethylbu...	206	C14H22O	000140-66-9	72
4			Hexestrol	270	C18H22O2	000084-16-2	72
5			Hexestrol, O-(n-propyloxycarbonyl)-	356	C22H28O4	1000487-65-1	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM011422\
Data File : BM033736.D
Acq On : 14 Jan 2022 20:00
Operator : CG/JU
Sample : N1121-03
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Instrument :
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ClientSampleId :
WASTE-CLASS-LAUNDRY-PIT

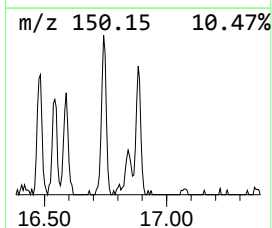
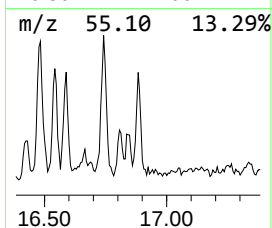
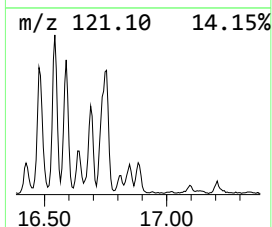
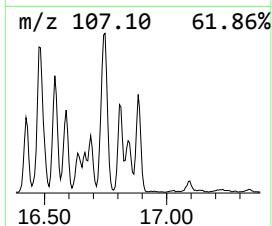
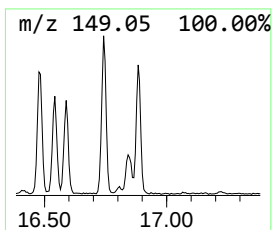
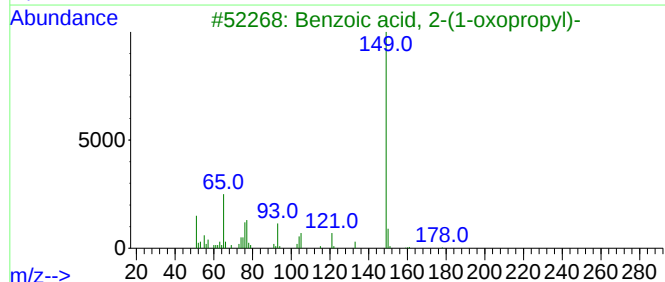
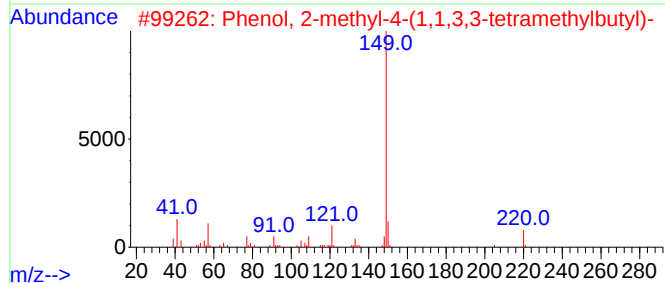
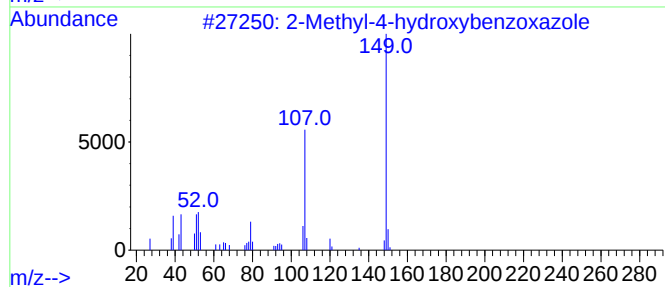
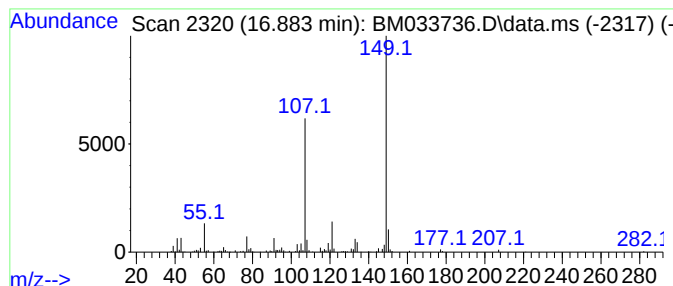
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 2-Methyl-4-hydroxybenzoxazole Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.883	2.29 ng	208646	Phenanthrene-d10	17.336

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Methyl-4-hydroxybenzoxazole	149	C8H7NO2	051110-60-2	64
2			Phenol, 2-methyl-4-(1,1,3,3-tetr...	220	C15H24O	002219-84-3	50
3			Benzoic acid, 2-(1-oxopropyl)-	178	C10H10O3	002360-45-4	50
4			6-Methylthieno[2,3-b]pyridine	149	C8H7NS	001759-30-4	49
5			7-Methylthieno[3,2-b]pyridine	149	C8H7NS	013362-83-9	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM011422\
Data File : BM033736.D
Acq On : 14 Jan 2022 20:00
Operator : CG/JU
Sample : N1121-03
Misc :
ALS Vial : 13 Sample Multiplier: 1

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ClientSampleId :
WASTE-CLASS-LAUNDRY-PIT

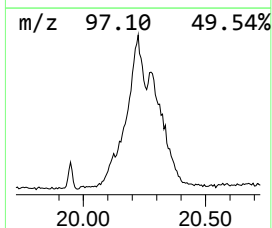
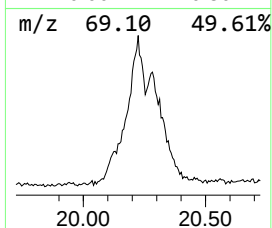
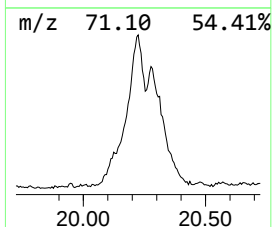
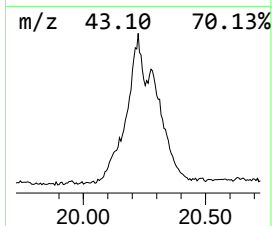
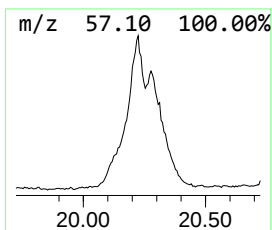
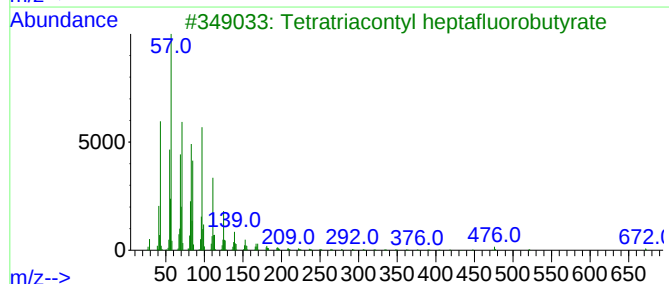
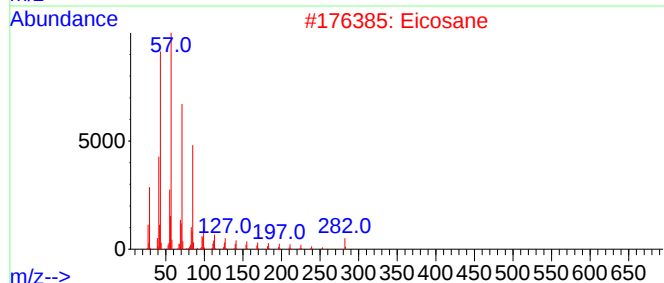
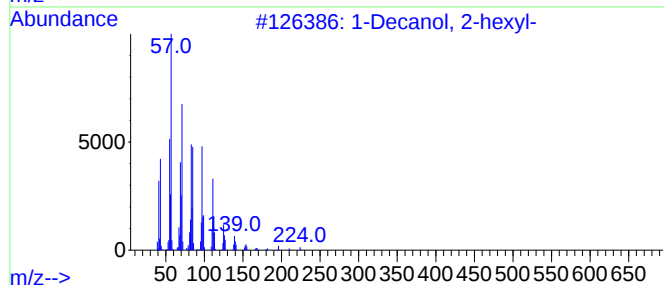
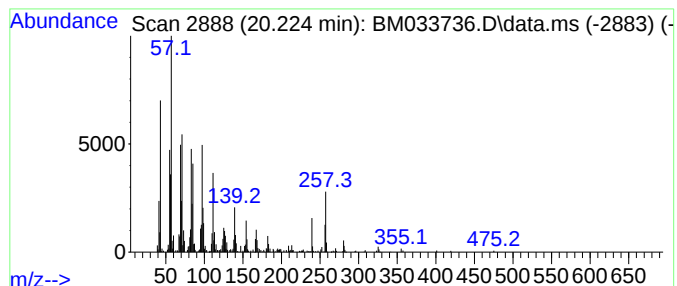
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 1-Decanol, 2-hexyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.224	6.49 ng	509443	Chrysene-d12	21.494

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Decanol, 2-hexyl-	242	C16H34O	002425-77-6	74
2			Eicosane	282	C20H42	000112-95-8	72
3			Tetratriacontyl heptafluorobutyrate	691	C38H69F7O2	1000351-84-1	68
4			1-Dodecanol, 2-octyl-	298	C20H42O	005333-42-6	68
5			Tetratriacontyl pentafluoropropi...	641	C37H69F5O2	1000351-81-5	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM011422\
Data File : BM033736.D
Acq On : 14 Jan 2022 20:00
Operator : CG/JU
Sample : N1121-03
Misc :
ALS Vial : 13 Sample Multiplier: 1

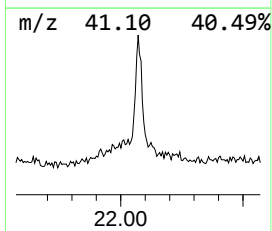
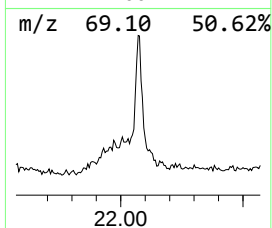
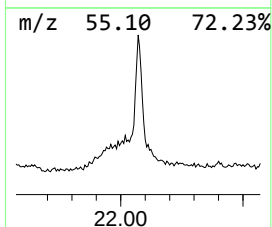
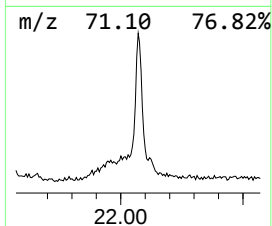
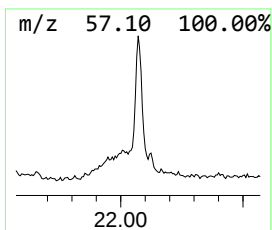
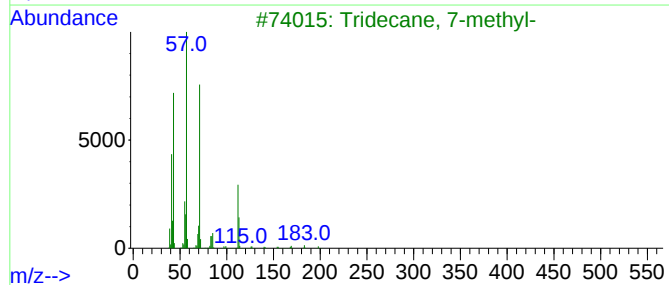
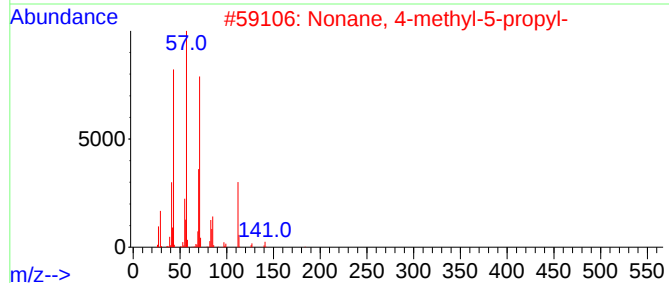
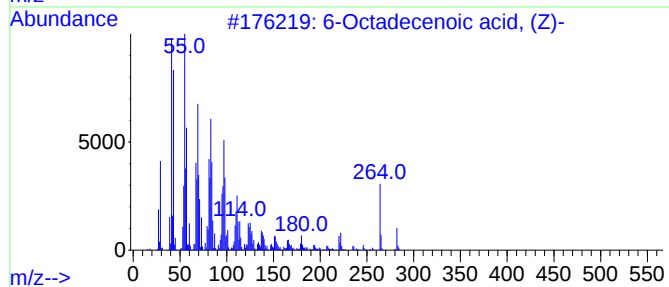
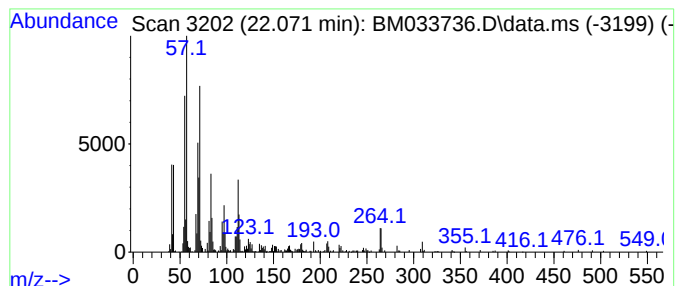
Instrument :
BNA_M
ClientSampleId :
WASTE-CLASS-LAUNDRY-PIT

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM010522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 6-Octadecenoic acid, (Z)- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
22.071	7.04 ng	553331	Chrysene-d12	21.494	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	6-Octadecenoic acid, (Z)-	282	C18H34O2	000593-39-5	90
2	Nonane, 4-methyl-5-propyl-	184	C13H28	062185-55-1	60
3	Tridecane, 7-methyl-	198	C14H30	026730-14-3	53
4	Carbonic acid, octyl vinyl ester	200	C11H20O3	1000383-25-5	50
5	(2E)-Dodec-2-en-1-yl methyl ether	198	C13H26O	1000334-06-3	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM011422\
Data File : BM033736.D
Acq On : 14 Jan 2022 20:00
Operator : CG/JU
Sample : N1121-03
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
WASTE-CLASS-LAUNDRY-PIT

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM010522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT 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
n-Propyl acetate	3.360	5.9	ng	297681	1	7.966	1002010	20.0
Propanoic acid,...	4.543	8.2	ng	413068	1	7.966	1002010	20.0
Isopropyl butyrate	5.013	5.2	ng	259913	1	7.966	1002010	20.0
unknown5.037	5.037	4.5	ng	224716	1	7.966	1002010	20.0
Phenol, 4-(1,1,...	16.424	3.9	ng	353861	4	17.336	1820080	20.0
5H-Pyrrolo(3,2-...	16.483	5.6	ng	508677	4	17.336	1820080	20.0
4-(7-Methylocty...	16.542	4.2	ng	380336	4	17.336	1820080	20.0
Phenol, 2-(1,1-...	16.589	2.5	ng	232213	4	17.336	1820080	20.0
1,3-Cyclopentad...	16.742	5.1	ng	464478	4	17.336	1820080	20.0
4-tert-Butylphe...	16.807	4.1	ng	375302	4	17.336	1820080	20.0
2-Methyl-4-hydr...	16.883	2.3	ng	208646	4	17.336	1820080	20.0
1-Decanol, 2-he...	20.224	6.5	ng	509443	5	21.494	1571030	20.0
6-Octadecenoic ...	22.071	7.0	ng	553331	5	21.494	1571030	20.0