

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112120\
 Data File : BP004044.D
 Acq On : 21 Nov 2020 18:39
 Operator : CG/JU
 Sample : L4839-05
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 B3-GW

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_P\METHODS\8270E-BP110320.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP004044.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.075	25	32	42	rBV	150301	303719	4.82%	0.975%
2	3.164	42	47	65	rVB	759909	1060630	16.83%	3.404%
3	5.263	400	404	413	rVB	86199	127100	2.02%	0.408%
4	5.805	488	496	512	rBV	1014611	1577860	25.03%	5.064%
5	6.399	590	597	604	rVB	49477	75209	1.19%	0.241%
6	7.452	769	776	791	rBV	555800	966873	15.34%	3.103%
7	8.263	907	914	923	rBV	390925	626131	9.93%	2.010%
8	9.428	1105	1112	1124	rBV	1379748	2287442	36.29%	7.341%
9	11.093	1387	1395	1414	rBV	497315	851787	13.51%	2.734%
10	13.510	1797	1806	1819	rBV	3184093	4641420	73.63%	14.896%
11	13.775	1846	1851	1858	rVB	63040	96626	1.53%	0.310%
12	14.310	1936	1942	1951	rBV	135537	195992	3.11%	0.629%
13	14.887	2033	2040	2049	rBV	732498	1071238	16.99%	3.438%
14	16.375	2285	2293	2308	rBV	2424386	3651296	57.92%	11.719%
15	16.698	2343	2348	2352	rBV	91803	126691	2.01%	0.407%
16	16.751	2352	2357	2364	rVV2	111755	208438	3.31%	0.669%
17	16.816	2364	2368	2372	rVV	102886	137579	2.18%	0.442%
18	16.863	2372	2376	2381	rVB2	71577	93231	1.48%	0.299%
19	17.016	2397	2402	2408	rVB4	64186	122649	1.95%	0.394%
20	17.081	2408	2413	2417	rBV	102701	138166	2.19%	0.443%
21	17.157	2423	2426	2432	rVB	42695	67750	1.07%	0.217%
22	17.510	2481	2486	2492	rBV	51360	79320	1.26%	0.255%
23	17.622	2498	2505	2510	rBV	873319	1274795	20.22%	4.091%
24	18.469	2644	2649	2663	rBV	620875	1123408	17.82%	3.606%
25	18.892	2717	2721	2729	rBV	112751	161132	2.56%	0.517%
26	18.963	2730	2733	2742	rVB	54901	78330	1.24%	0.251%
27	19.674	2850	2854	2865	rBV	170609	288075	4.57%	0.925%
28	20.116	2924	2929	2935	rVB	5144892	6303721	100.00%	20.231%
29	20.674	3021	3024	3029	rVB	101101	118047	1.87%	0.379%
30	21.310	3128	3132	3142	rBV	309605	547615	8.69%	1.758%
31	21.651	3185	3190	3196	rBV	1046730	1365598	21.66%	4.383%
32	24.115	3603	3609	3621	rVB2	657220	1390176	22.05%	4.462%

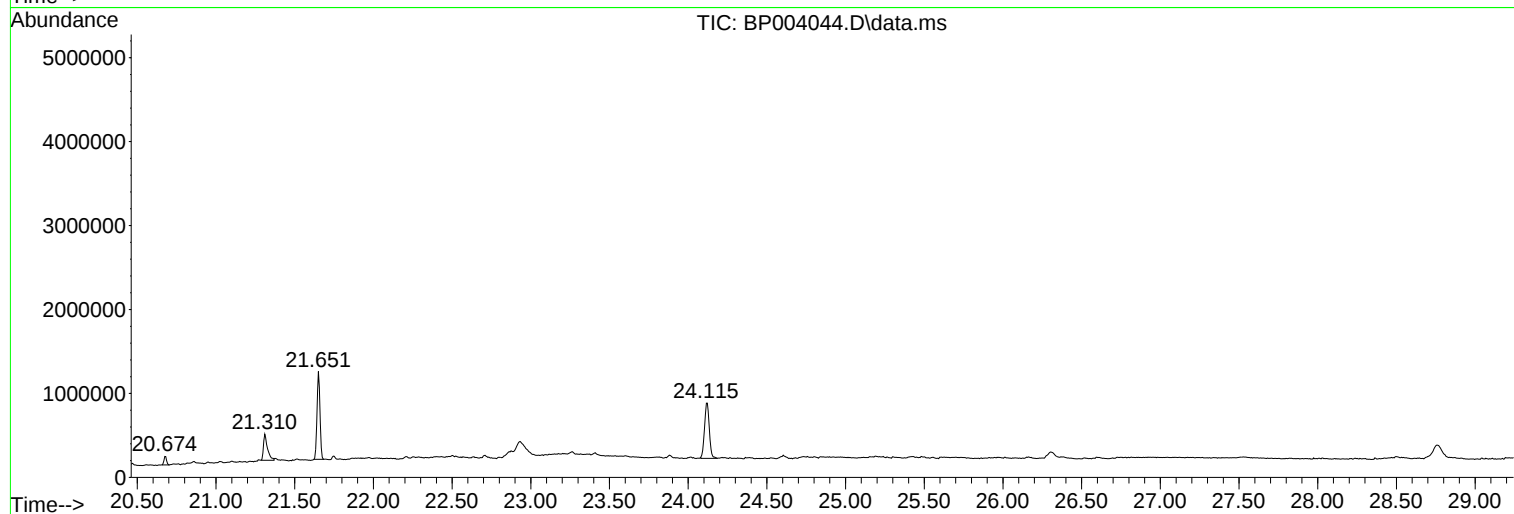
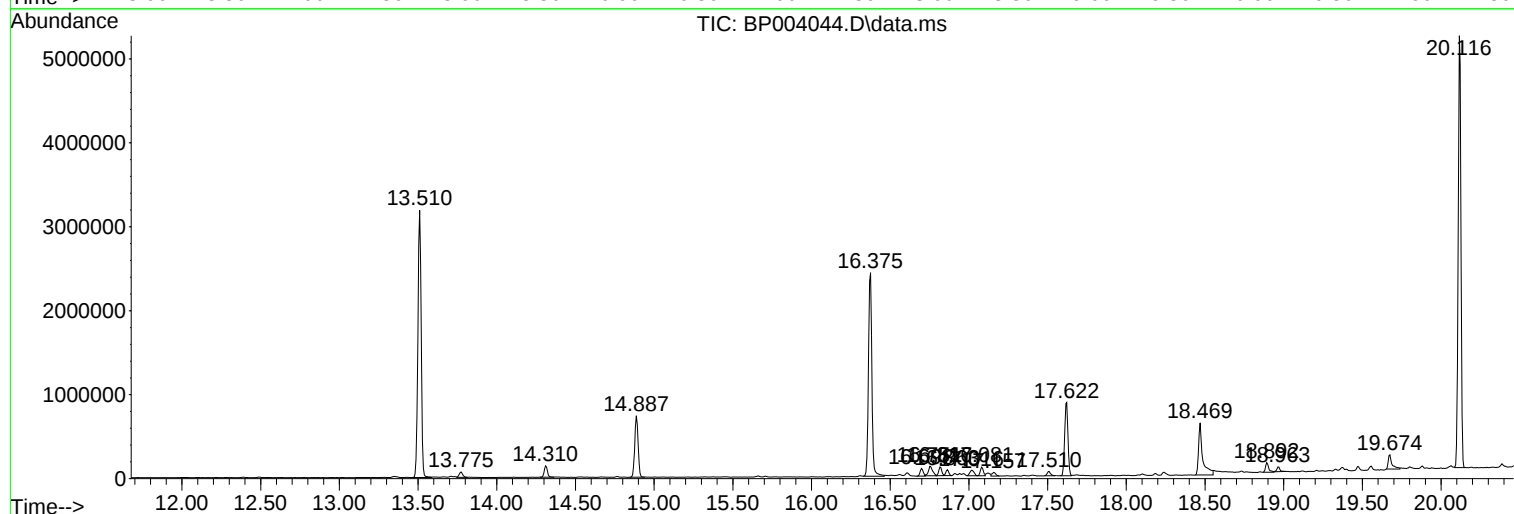
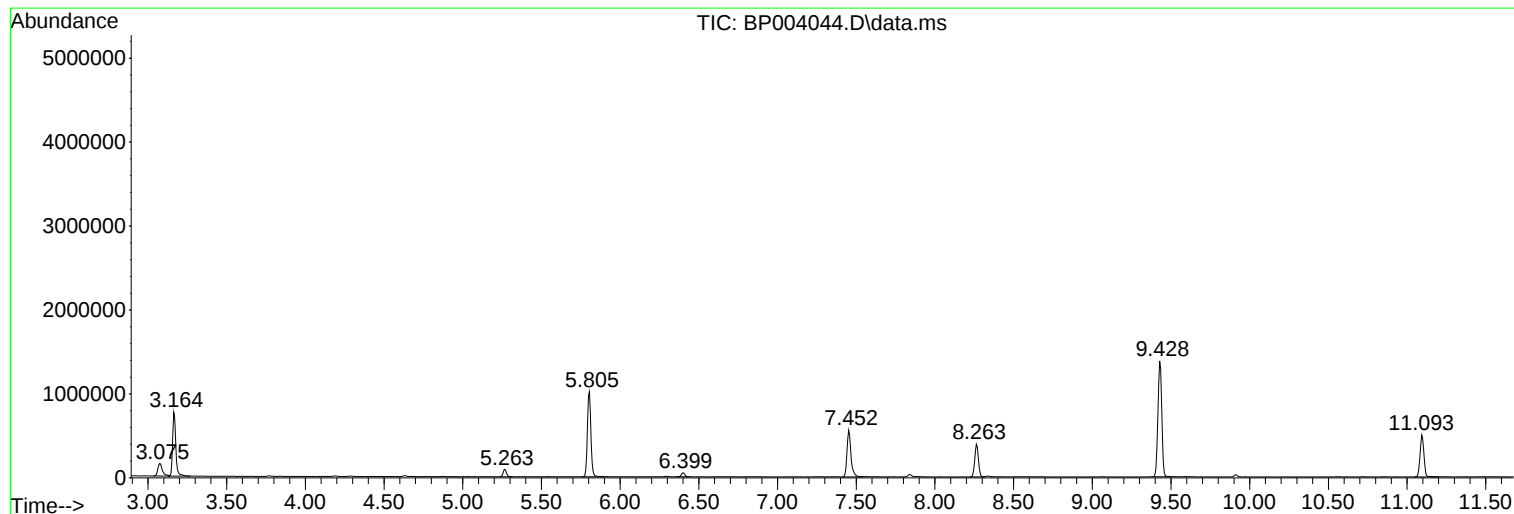
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270E-BP110320.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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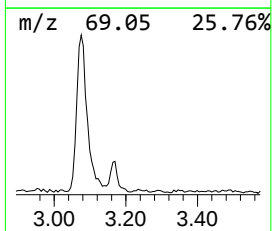
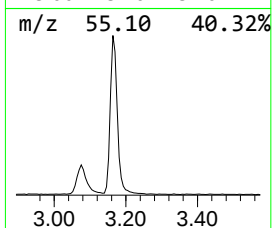
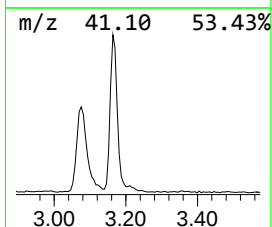
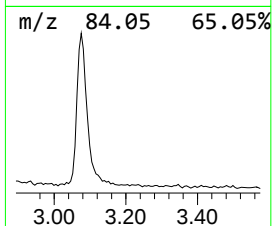
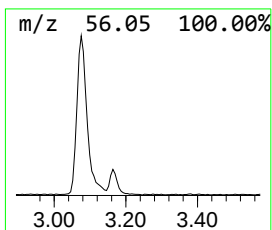
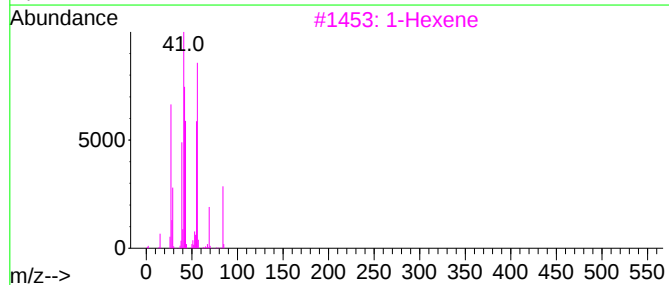
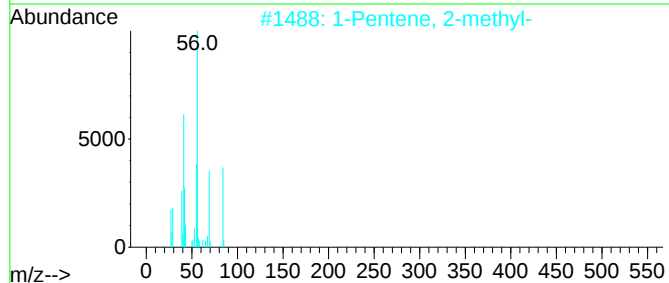
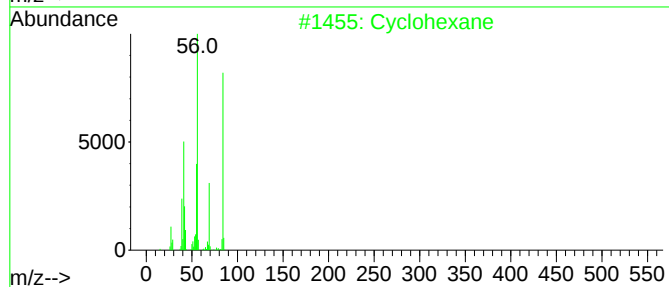
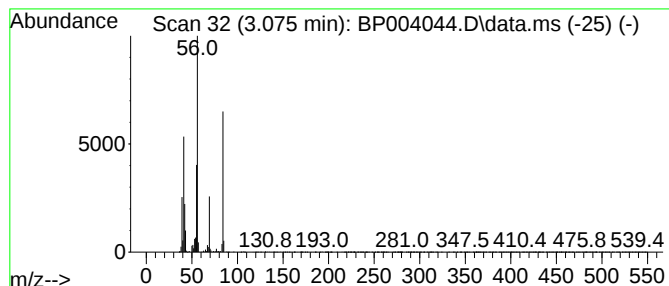
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270E-BP110320.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Cyclohexane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.075	9.70 ng	303719	1,4-Dichlorobenzene-d4	8.263

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclohexane	84	C6H12	000110-82-7	95
2		1-Pentene, 2-methyl-	84	C6H12	000763-29-1	72
3		1-Hexene	84	C6H12	000592-41-6	59
4		Cyclobutane	56	C4H8	000287-23-0	49
5		1-Hexene, 5-methyl-	98	C7H14	003524-73-0	47



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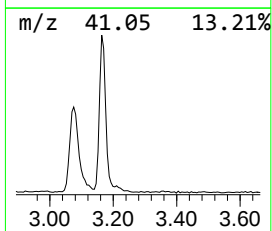
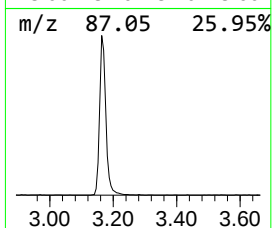
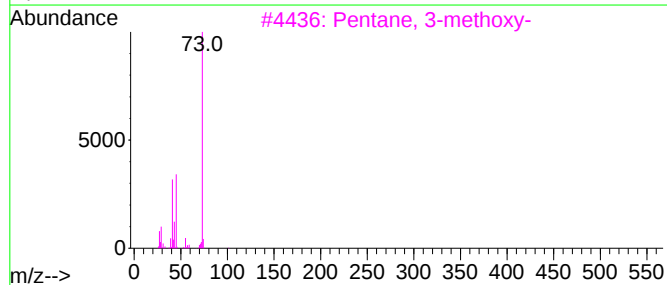
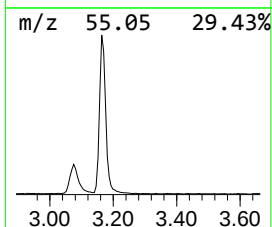
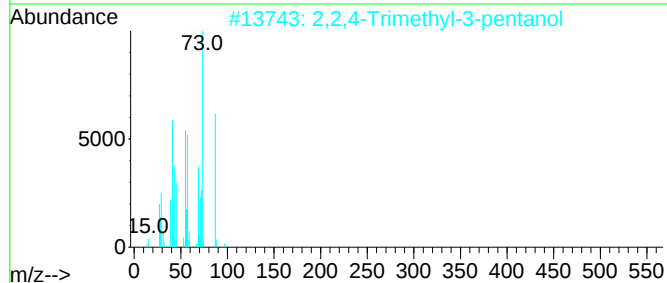
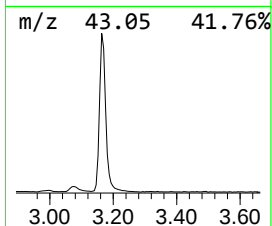
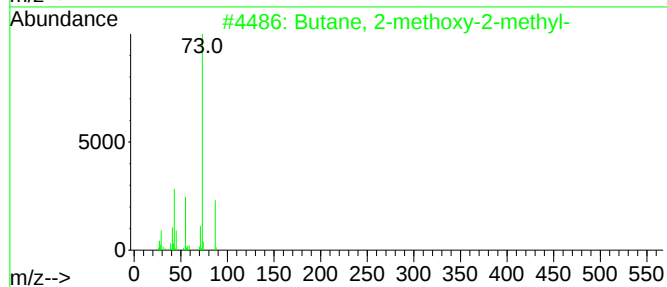
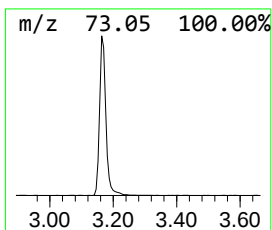
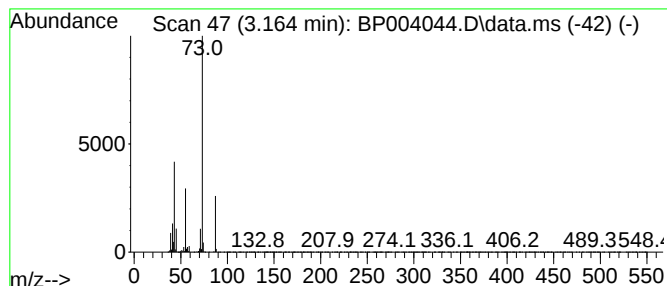
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270E-BP110320.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Butane, 2-methoxy-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.164	33.88 ng	1060630	1,4-Dichlorobenzene-d4	8.263

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
3			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12
4			2-Methyl-5-hexen-3-ol	114	C7H14O	032815-70-6	12
5			N-Ethylformamide	73	C3H7NO	000627-45-2	9



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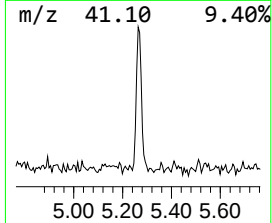
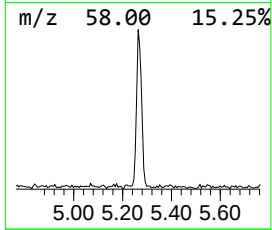
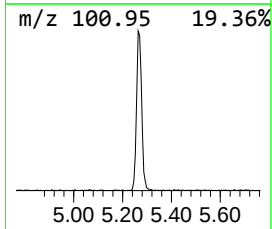
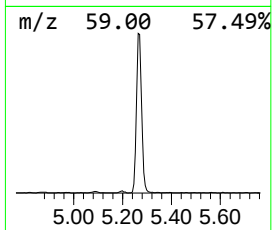
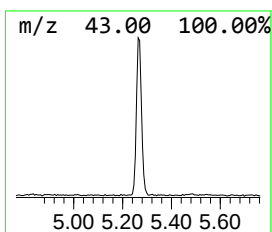
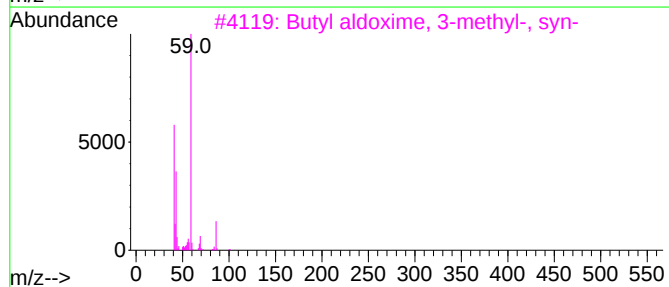
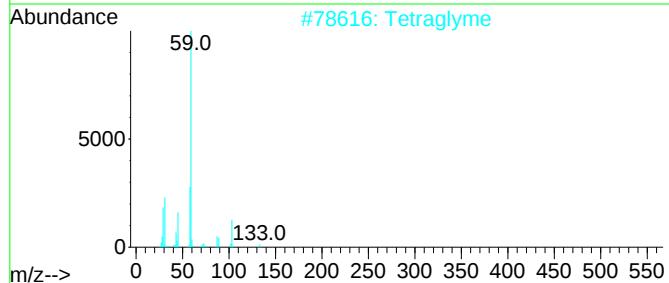
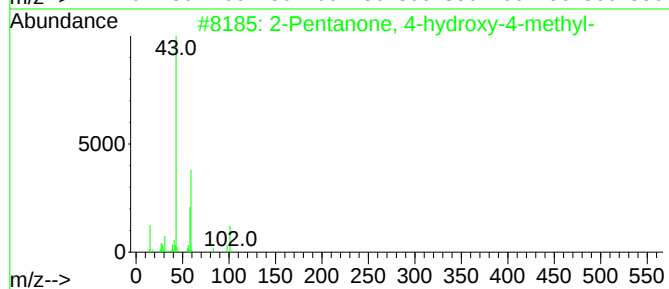
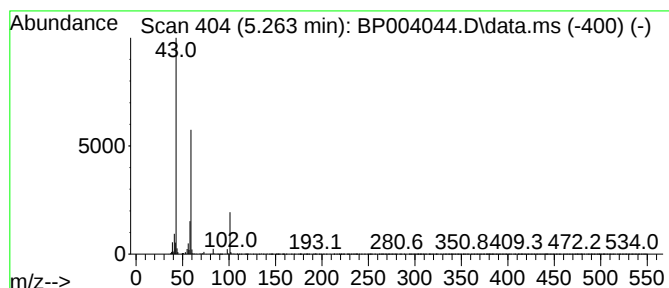
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD		R.T.	
5.263	4.06 ng	127100	1,4-Dichlorobenzene-d4		8.263	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-		116	C6H12O2	000123-42-2	56
2	Tetraglyme		222	C10H22O5	000143-24-8	25
3	Butyl aldoxime, 3-methyl-, syn-		101	C5H11NO	005780-40-5	25
4	2,3-Butanedione, monooxime		101	C4H7NO2	000057-71-6	25
5	2,5,8,11-Tetraoxadodecane		178	C8H18O4	000112-49-2	23



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ClientSampleId :
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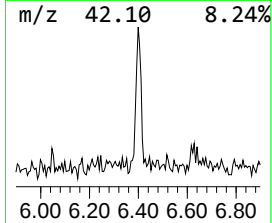
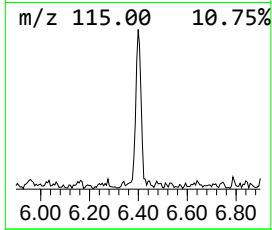
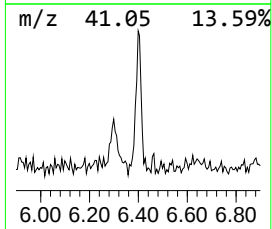
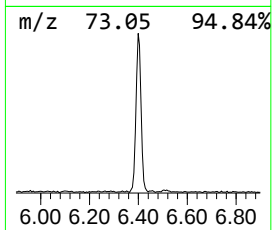
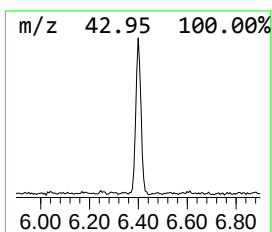
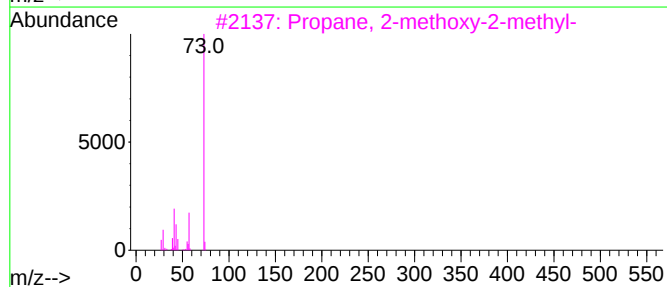
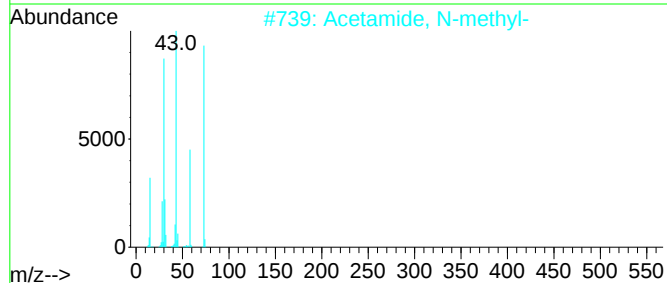
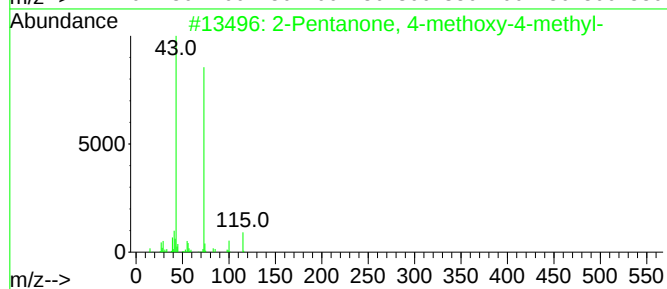
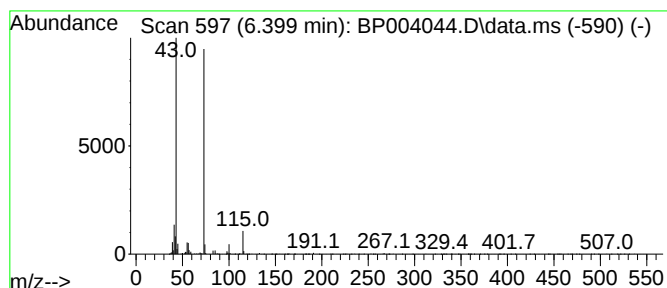
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Acetamide, N-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.399	2.40 ng	75209	1,4-Dichlorobenzene-d4	8.263

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-methoxy-4-methyl-	130	C7H14O2	000107-70-0	78		
2	Acetamide, N-methyl-	73	C3H7NO	000079-16-3	52		
3	Propane, 2-methoxy-2-methyl-	88	C5H12O	001634-04-4	12		
4	1,3-Dioxolane, 2-(1-methylethyl)-	116	C6H12O2	000822-83-3	9		
5	2-Propyn-1-ol, acetate	98	C5H6O2	000627-09-8	9		



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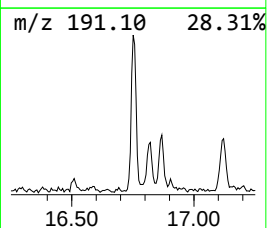
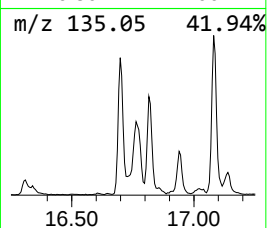
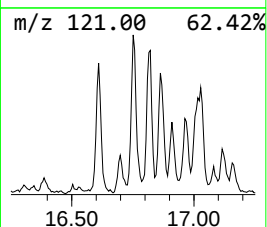
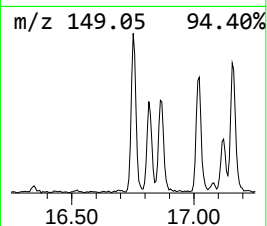
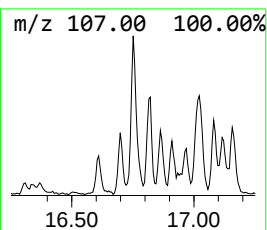
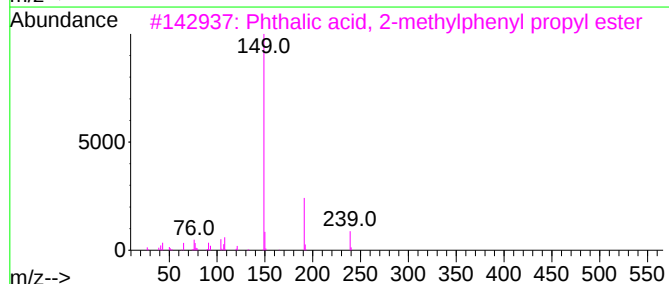
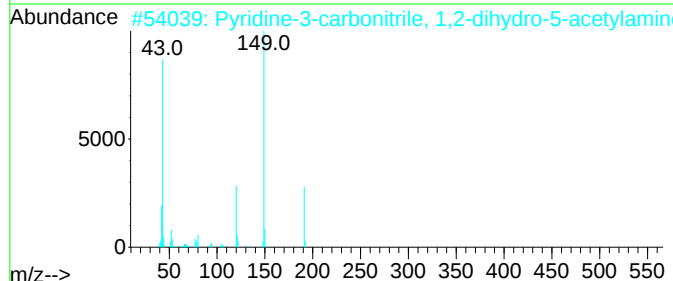
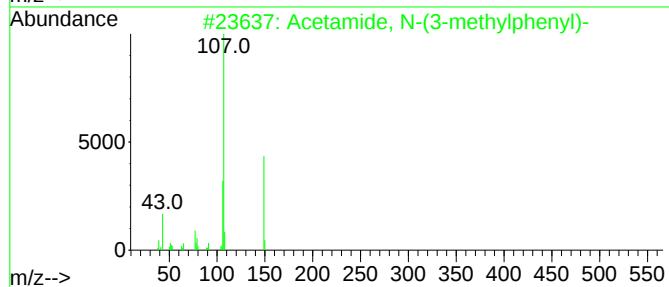
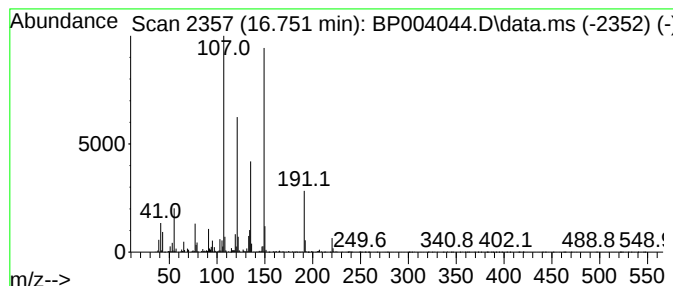
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 unknown16.751 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.751	3.27 ng	208438	Phenanthrene-d10	17.622

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acetamide, N-(3-methylphenyl)-	149	C9H11NO	000537-92-8	38
2			Pyridine-3-carbonitrile, 1,2-dih...	191	C9H9N3O2	220098-92-0	38
3			Phthalic acid, 2-methylphenyl pr...	298	C18H18O4	1000314-95-1	35
4			Phthalic acid, 3-methylphenyl pr...	298	C18H18O4	1000315-36-6	35
5			Phthalic acid, phenyl propyl ester	284	C17H16O4	1000314-86-9	35



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ClientSampleId :
B3-GW

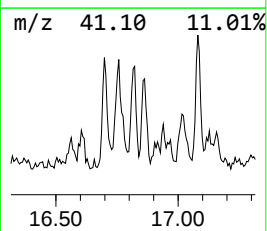
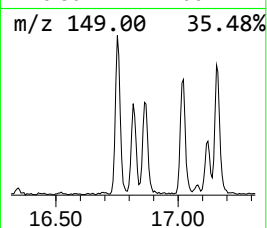
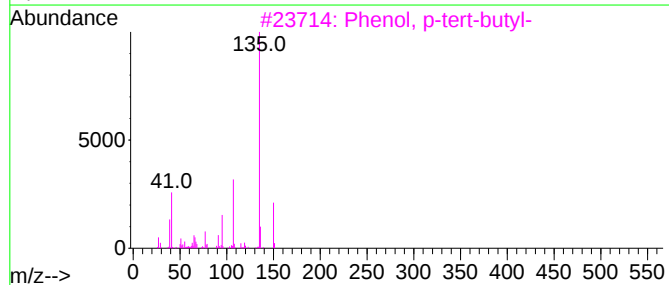
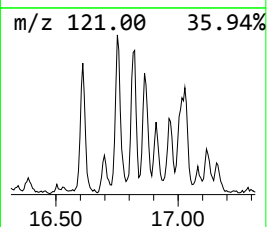
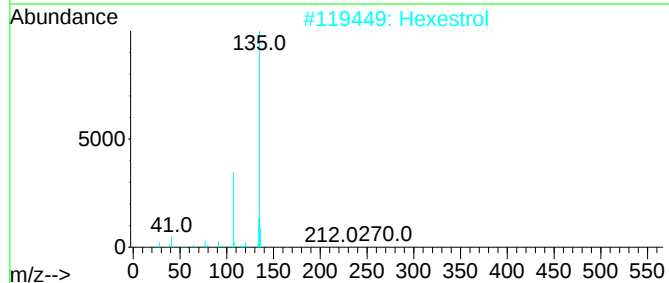
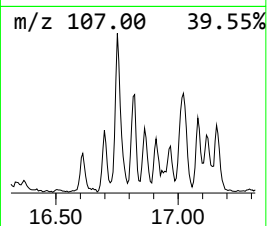
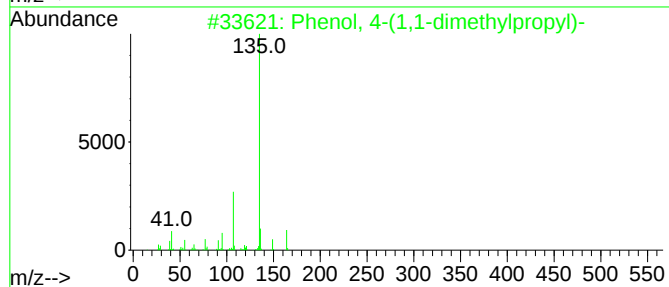
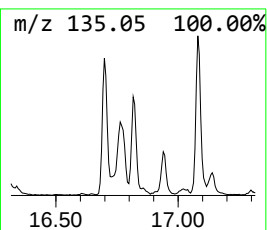
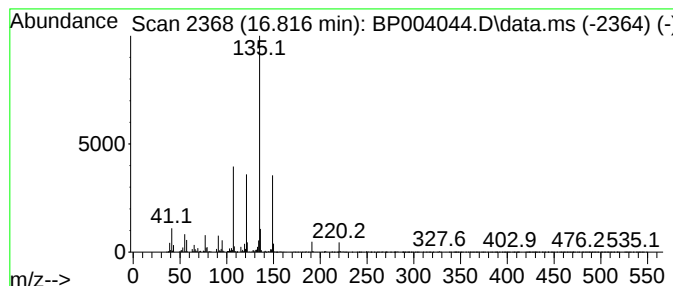
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Phenol, 4-(1,1-dimethylprop... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.816	2.16 ng	137579	Phenanthrene-d10	17.622

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	59
2			Hexestrol	270	C18H22O2	000084-16-2	53
3			Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	53
4			Adamantane-1-carboxylic acid, 4-...	284	C18H20O3	1000276-61-1	46
5			N-(2-Ethyl-3-hydroxy-6-methyl-4-...	314	C19H26N2O2	1000260-99-4	45



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112120\
Data File : BP004044.D
Acq On : 21 Nov 2020 18:39
Operator : CG/JU
Sample : L4839-05
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
B3-GW

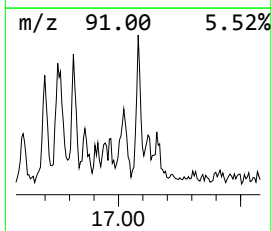
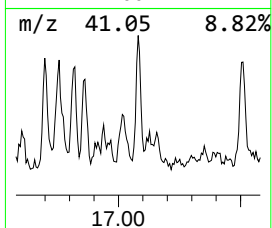
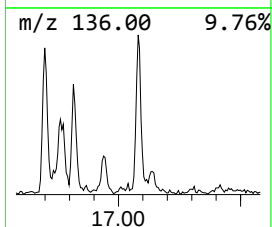
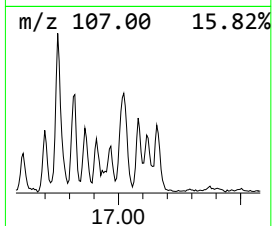
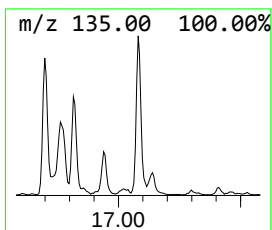
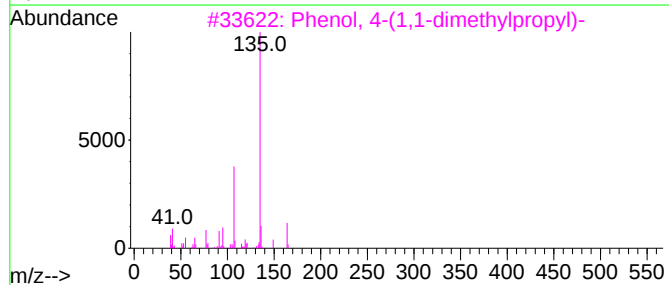
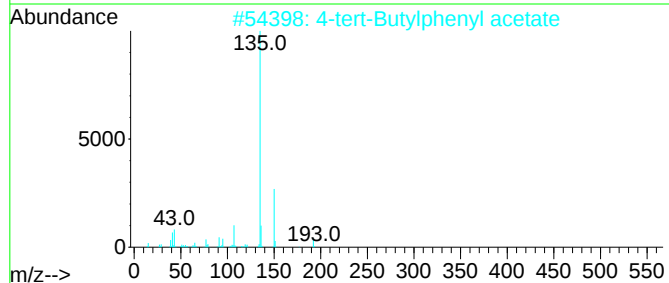
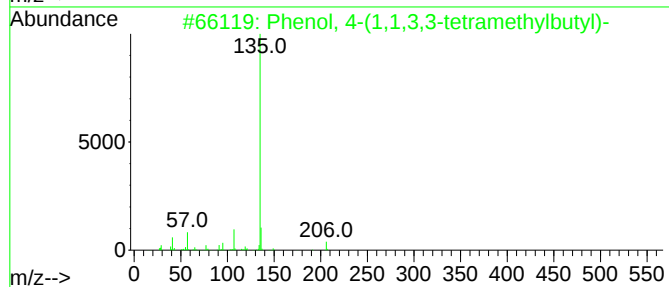
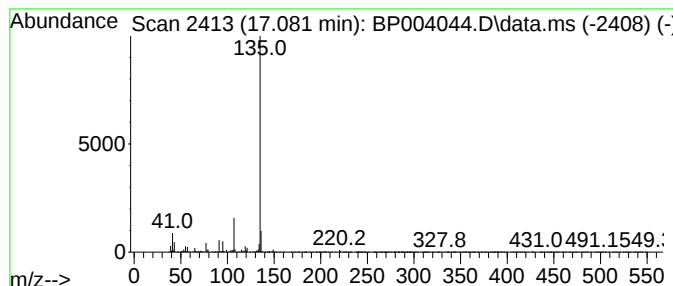
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Phenol, 4-(1,1,3,3-tetramet... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.081	2.17 ng	138166	Phenanthrene-d10	17.622

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 4-(1,1,3,3-tetramethylbu...	206	C14H22O	000140-66-9	78
2			4-tert-Butylphenyl acetate	192	C12H16O2	003056-64-2	78
3			Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	78
4			p-Methoxybenzoic acid, 2-methylp...	242	C15H14O3	007504-73-6	64
5			Hexestrol	270	C18H22O2	000084-16-2	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112120\
Data File : BP004044.D
Acq On : 21 Nov 2020 18:39
Operator : CG/JU
Sample : L4839-05
Misc :
ALS Vial : 8 Sample Multiplier: 1

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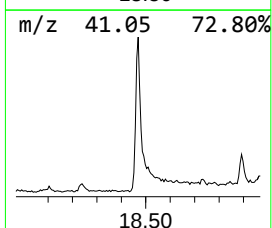
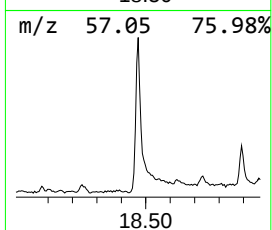
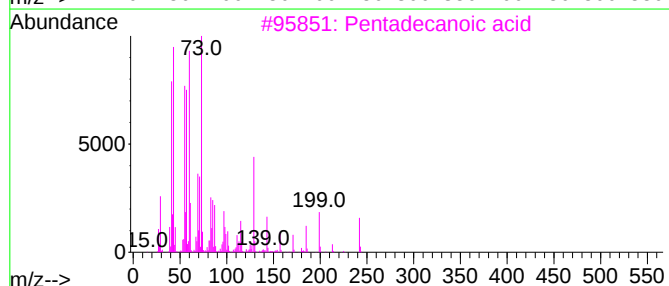
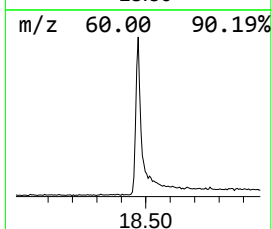
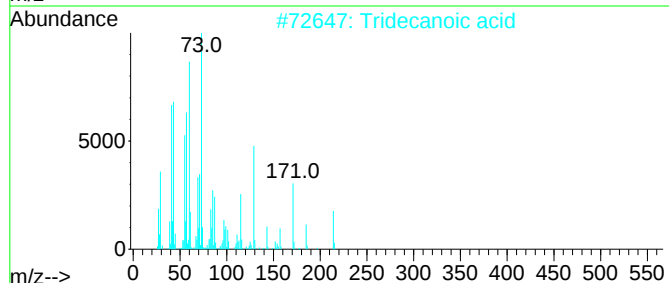
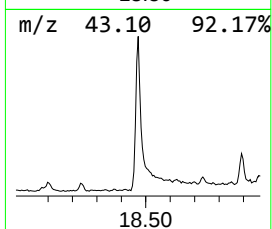
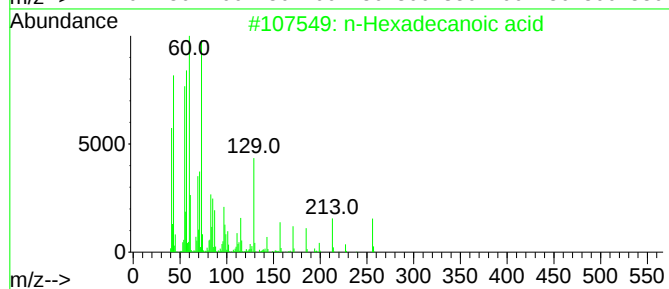
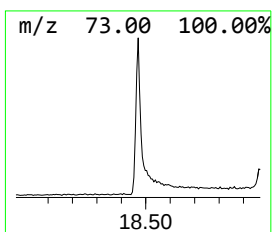
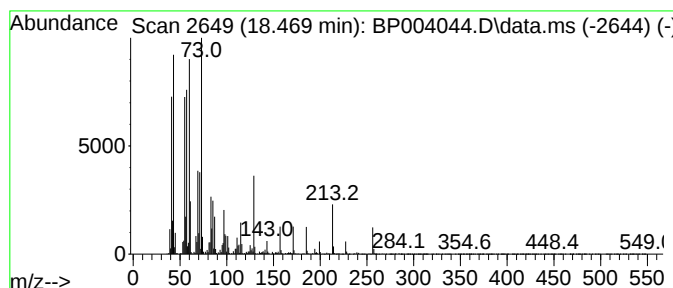
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.469	17.62 ng	1123410	Phenanthrene-d10	17.622

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		Tridecanoic acid	214	C13H26O2	000638-53-9	93
3		Pentadecanoic acid	242	C15H30O2	001002-84-2	87
4		Tetradecanoic acid	228	C14H28O2	000544-63-8	64
5		Octadecanoic acid	284	C18H36O2	000057-11-4	60



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112120\
Data File : BP004044.D
Acq On : 21 Nov 2020 18:39
Operator : CG/JU
Sample : L4839-05
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
B3-GW

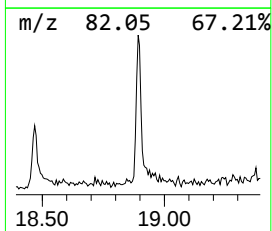
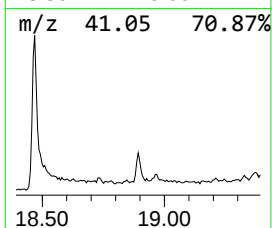
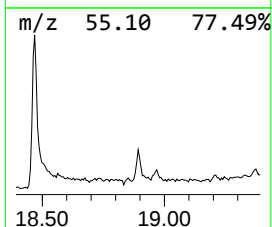
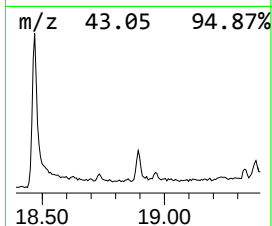
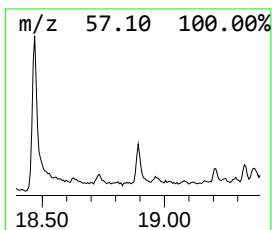
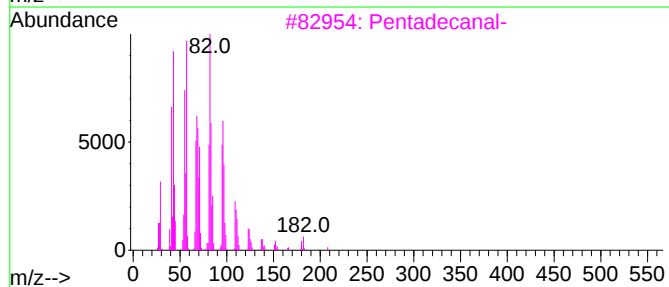
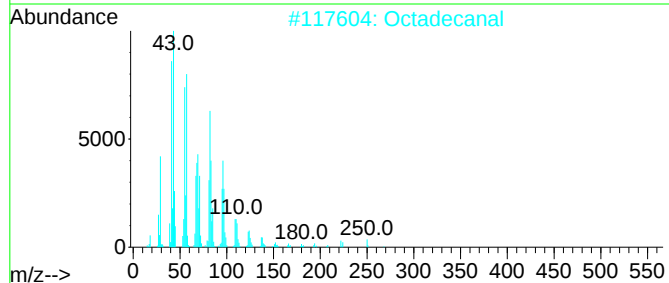
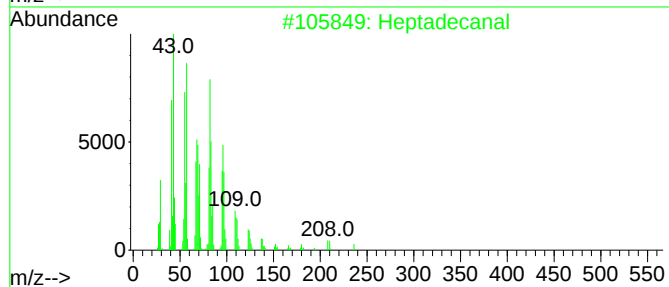
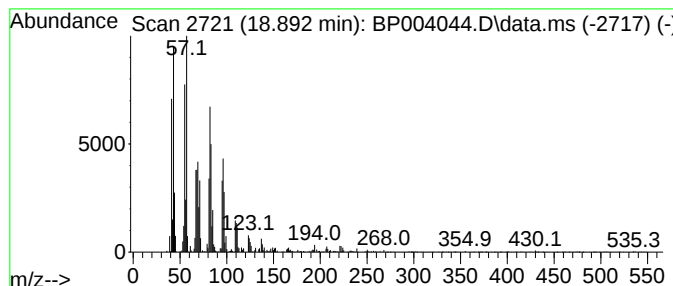
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Heptadecanal Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.892	2.53 ng	161132	Phenanthrene-d10	17.622

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecanal	254	C17H34O	1000376-70-0	96
2		Octadecanal	268	C18H36O	000638-66-4	94
3		Pentadecanal-	226	C15H30O	002765-11-9	94
4		Hexadecanal	240	C16H32O	000629-80-1	91
5		Z-2-Octadecen-1-ol	268	C18H36O	1000131-11-0	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112120\
Data File : BP004044.D
Acq On : 21 Nov 2020 18:39
Operator : CG/JU
Sample : L4839-05
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
B3-GW

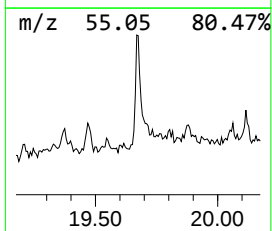
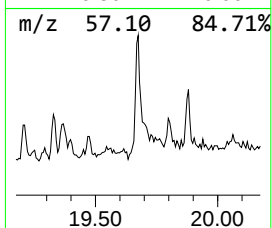
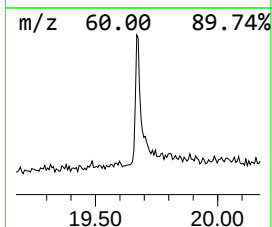
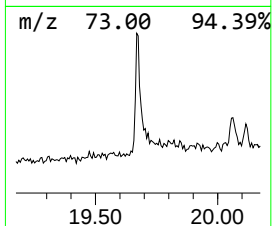
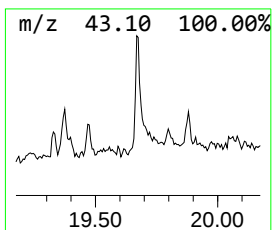
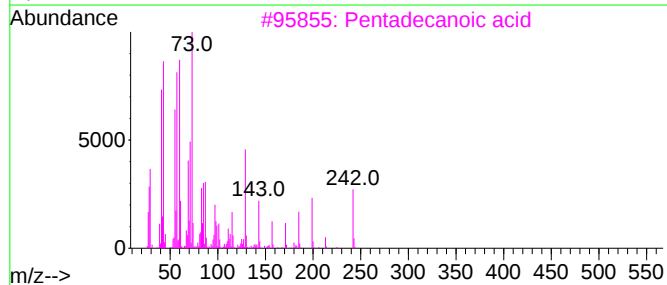
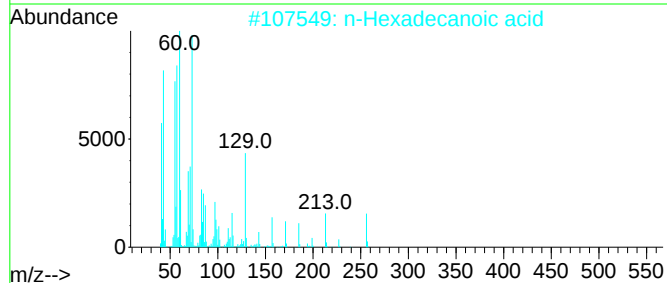
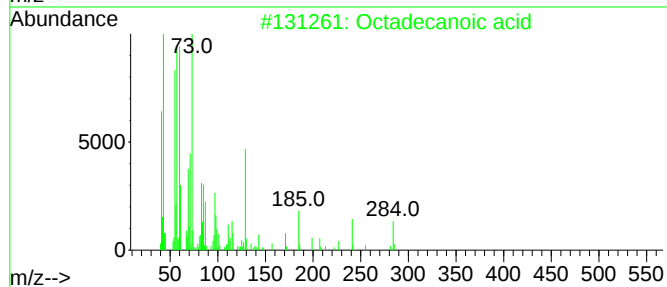
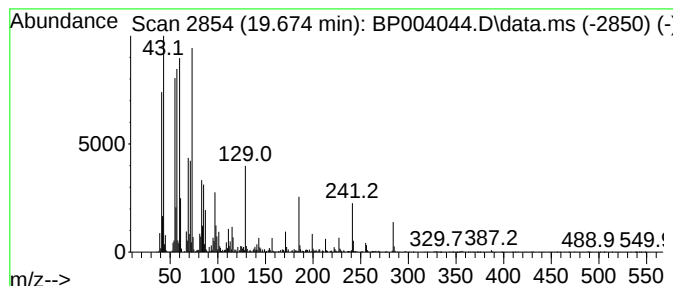
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Octadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.674	4.22 ng	288075	Chrysene-d12	21.651

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	99
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	91
3			Pentadecanoic acid	242	C15H30O2	001002-84-2	89
4			Tetradecanoic acid	228	C14H28O2	000544-63-8	83
5			Tridecanoic acid	214	C13H26O2	000638-53-9	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112120\
Data File : BP004044.D
Acq On : 21 Nov 2020 18:39
Operator : CG/JU
Sample : L4839-05
Misc :
ALS Vial : 8 Sample Multiplier: 1

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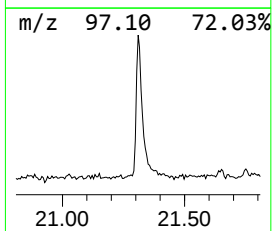
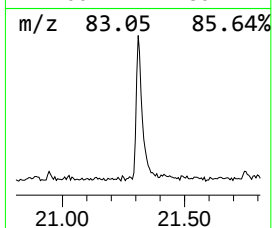
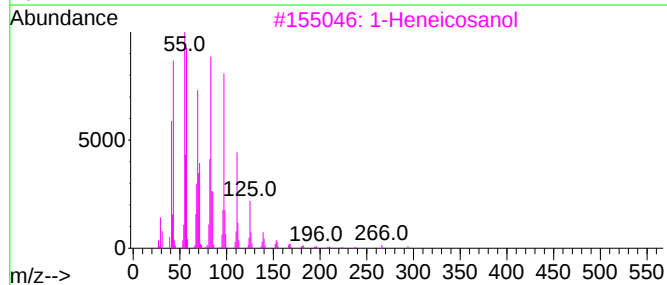
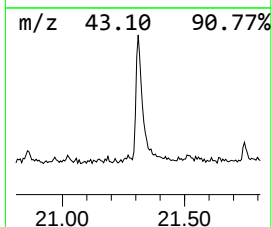
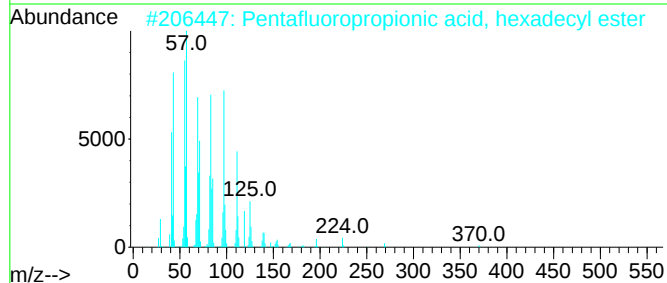
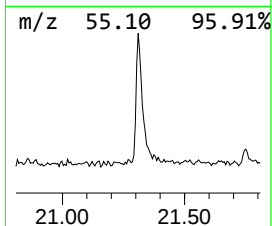
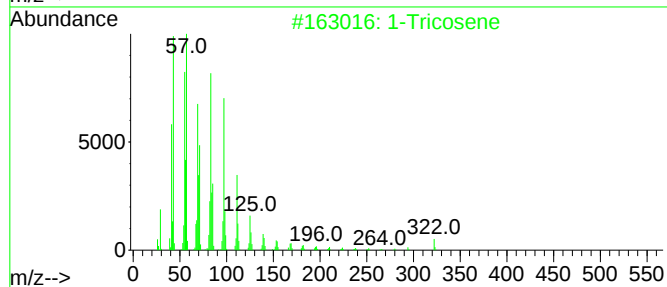
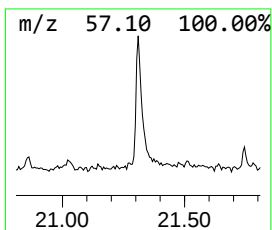
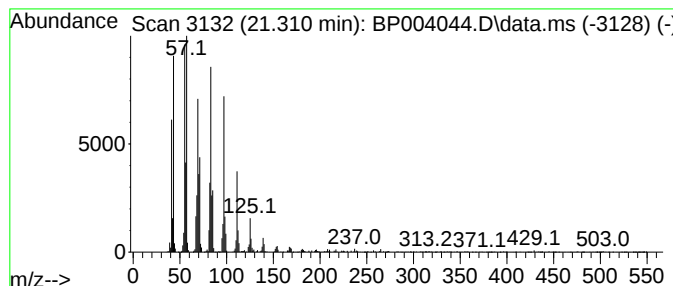
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 1-Tricosene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.310	8.02 ng	547615	Chrysene-d12	21.651

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Tricosene	322	C23H46	018835-32-0	98
2			Pentafluoropropionic acid, hexad...	388	C19H33F5O2	006222-07-7	94
3			1-Heneicosanol	312	C21H44O	015594-90-8	94
4			1-Docosene	308	C22H44	001599-67-3	94
5			Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112120\
Data File : BP004044.D
Acq On : 21 Nov 2020 18:39
Operator : CG/JU
Sample : L4839-05
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270E-BP110320.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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Butane, 2-metho...	3.164	33.9	ng	1060630	1	8.263	626131	20.0
2-Pentanone, 4-...	5.263	4.1	ng	127100	1	8.263	626131	20.0
Acetamide, N-me...	6.399	2.4	ng	75209	1	8.263	626131	20.0
unknown16.751	16.751	3.3	ng	208438	4	17.622	1274800	20.0
Phenol, 4-(1,1-...	16.816	2.2	ng	137579	4	17.622	1274800	20.0
Phenol, 4-(1,1,...	17.081	2.2	ng	138166	4	17.622	1274800	20.0
n-Hexadecanoic ...	18.469	17.6	ng	1123410	4	17.622	1274800	20.0
Heptadecanal	18.892	2.5	ng	161132	4	17.622	1274800	20.0
Octadecanoic acid	19.674	4.2	ng	288075	5	21.651	1365600	20.0
1-Tricosene	21.310	8.0	ng	547615	5	21.651	1365600	20.0