

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF053119\
 Data File : BF114733.D
 Acq On : 1 Jun 2019 3:13
 Operator : HP/JU
 Sample : K3072-03
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-3

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0
 Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF052919.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.316	542	549	552	rBV	7703859	9340244	94.82%	10.321%
2	6.340	716	723	726	rBV	7820207	9464970	96.09%	10.459%
3	6.469	739	745	748	rBV	9074747	9850559	100.00%	10.885%
4	6.698	780	784	787	rBV	2679187	2351359	23.87%	2.598%
5	6.857	806	811	814	rBV	7527119	7503823	76.18%	8.292%
6	7.257	873	879	882	rBV	5683398	6321406	64.17%	6.985%
7	7.975	996	1001	1004	rBV	3465821	2855650	28.99%	3.156%
8	9.051	1178	1184	1187	rBV	9460233	9351402	94.93%	10.334%
9	9.439	1246	1250	1257	rBV	1703102	1361326	13.82%	1.504%
10	9.722	1293	1298	1304	rBV	3414534	2983756	30.29%	3.297%
11	10.504	1426	1431	1434	rBV	5671910	5432630	55.15%	6.003%
12	11.192	1544	1548	1552	rBV	2989575	2564506	26.03%	2.834%
13	11.739	1637	1641	1649	rBV	1377131	1532951	15.56%	1.694%
14	12.522	1770	1774	1783	rBV	608749	817495	8.30%	0.903%
15	12.774	1813	1817	1821	rBV	6940787	7073271	71.81%	7.816%
16	13.680	1968	1971	1977	rVB	1423031	1400184	14.21%	1.547%
17	13.816	1990	1994	1997	rBV	2279934	2058574	20.90%	2.275%
18	14.010	2024	2027	2029	rBV	373992	344984	3.50%	0.381%
19	14.310	2075	2078	2082	rBV	603116	629381	6.39%	0.695%
20	14.345	2082	2084	2090	rVB	252589	310740	3.15%	0.343%
21	14.604	2125	2128	2134	rVB	854115	861411	8.74%	0.952%
22	14.915	2178	2181	2186	rVB	827932	859836	8.73%	0.950%
23	15.204	2224	2230	2235	rVB	1558771	1995399	20.26%	2.205%
24	15.263	2237	2240	2249	rVB	703859	915523	9.29%	1.012%
25	15.662	2304	2308	2312	rBV	458558	654133	6.64%	0.723%
26	15.704	2312	2315	2323	rVB	243446	417872	4.24%	0.462%
27	15.868	2339	2343	2354	rVB2	241154	560064	5.69%	0.619%
28	16.121	2382	2386	2390	rBV2	207648	312469	3.17%	0.345%
29	16.262	2406	2410	2418	rVB	213848	369949	3.76%	0.409%

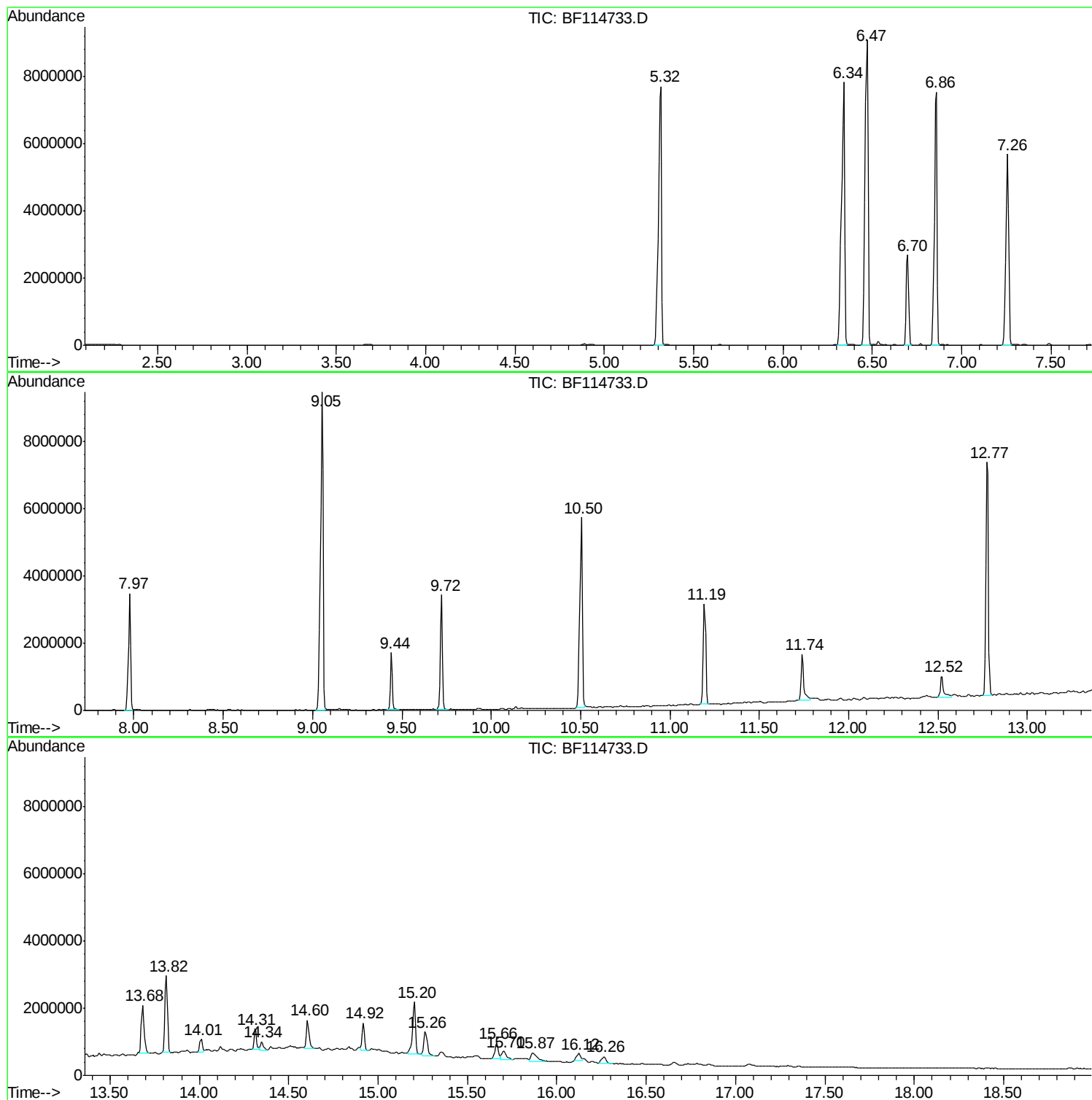
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF052919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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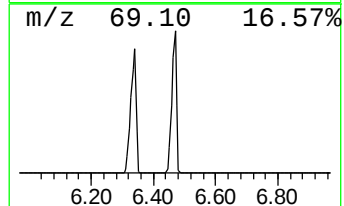
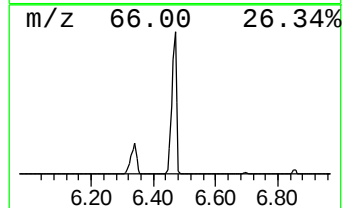
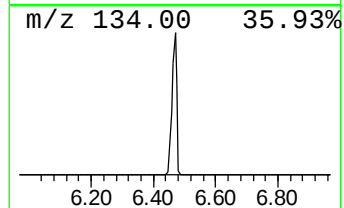
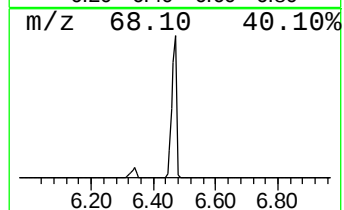
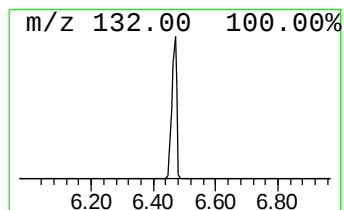
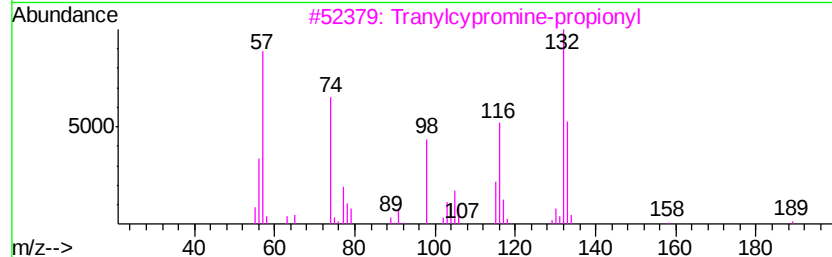
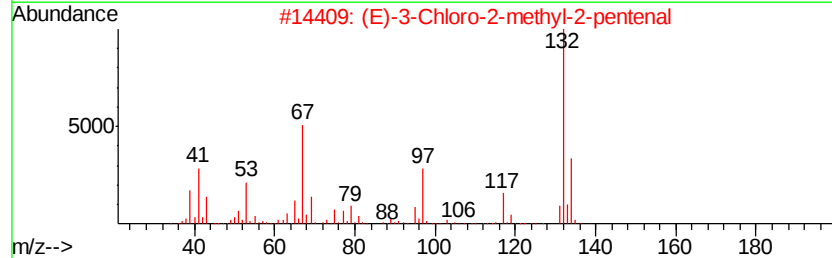
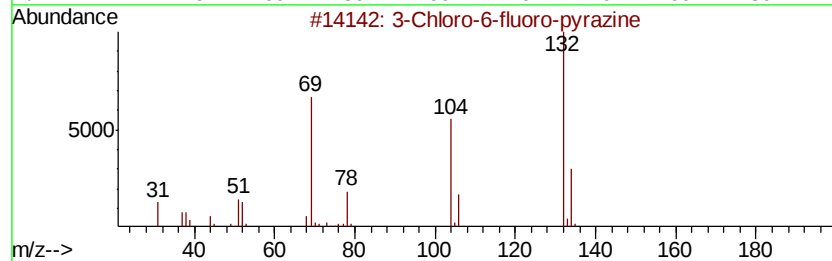
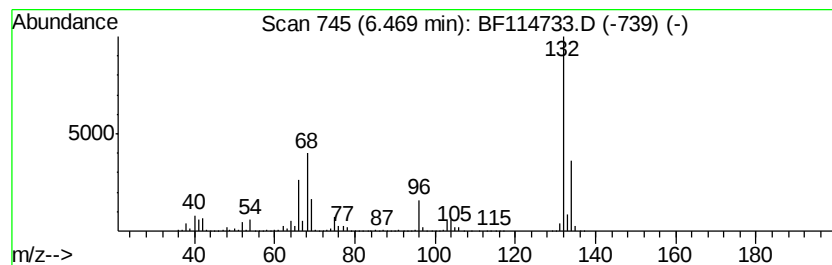
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown6.47 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.47	83.79 ng	9850560	1,4-Dichlorobenzene-d4	6.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
3		Tranvlcyproline-propionyl	189	C12H15NO	1000123-86-3	12
4		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9
5		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9



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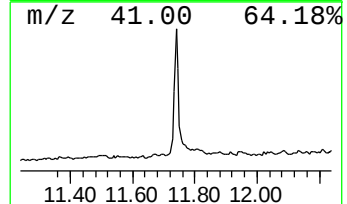
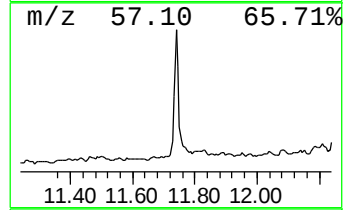
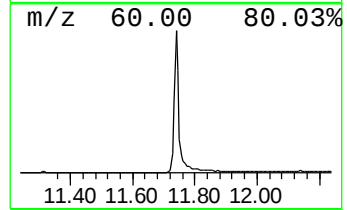
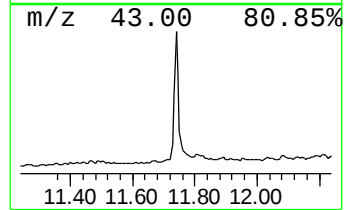
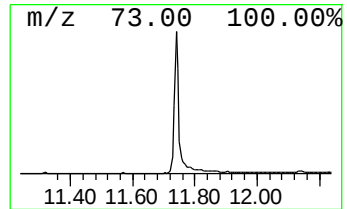
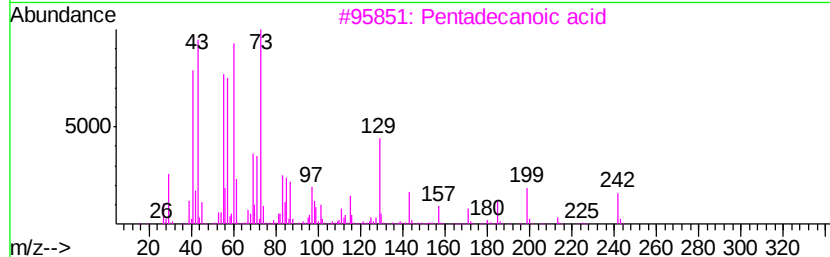
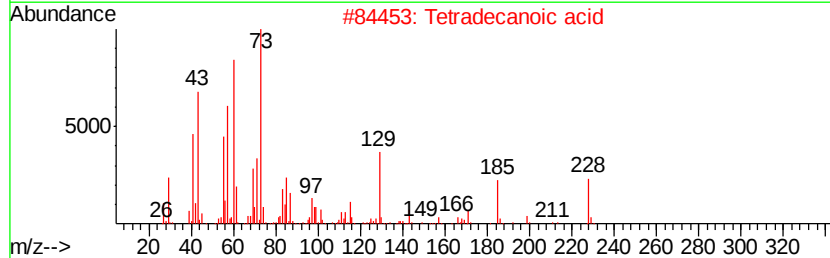
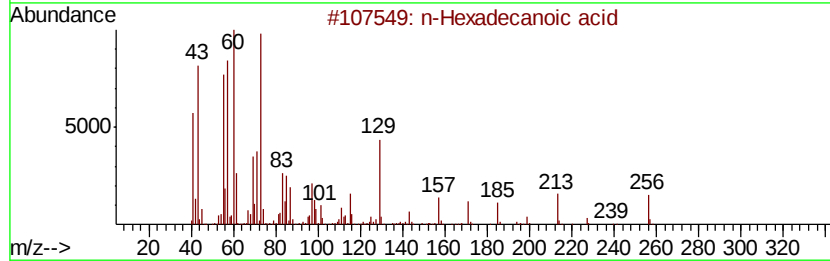
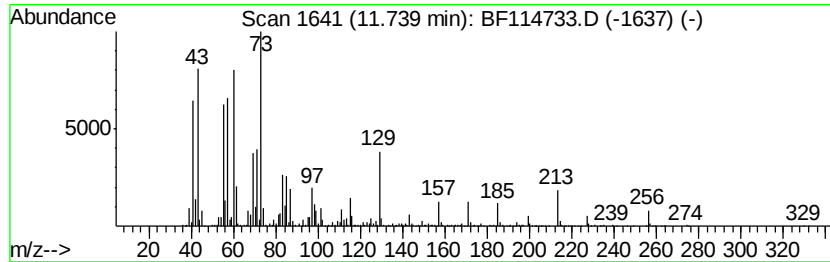
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 3 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.74	11.96 ng	1532950	Phenanthrene-d10	11.19

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		Tetradecanoic acid	228	C14H28O2	000544-63-8	94
3		Pentadecanoic acid	242	C15H30O2	001002-84-2	94
4		Tridecanoic acid	214	C13H26O2	000638-53-9	93
5		Octadecanoic acid	284	C18H36O2	000057-11-4	90



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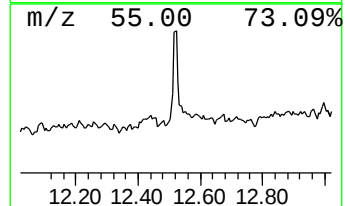
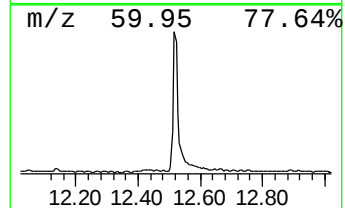
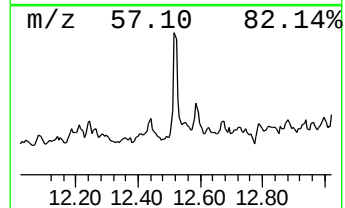
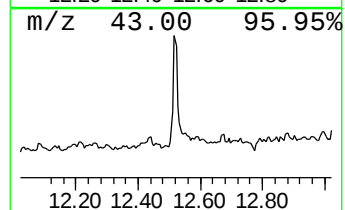
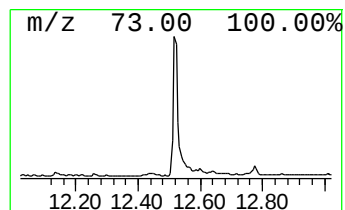
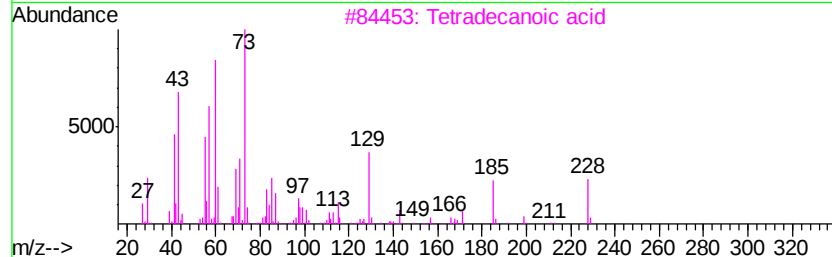
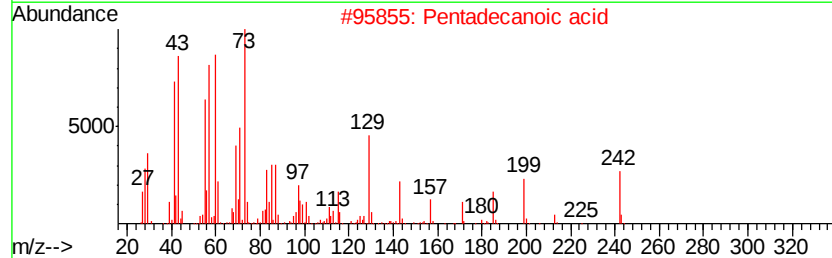
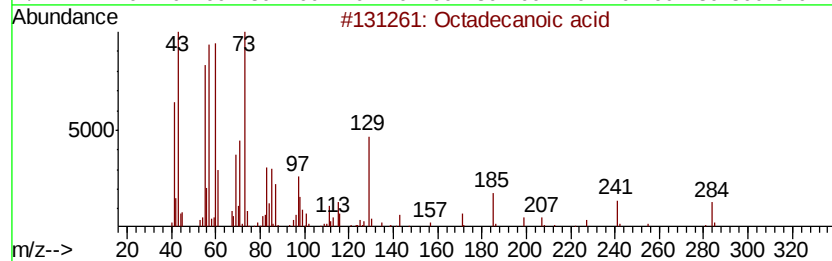
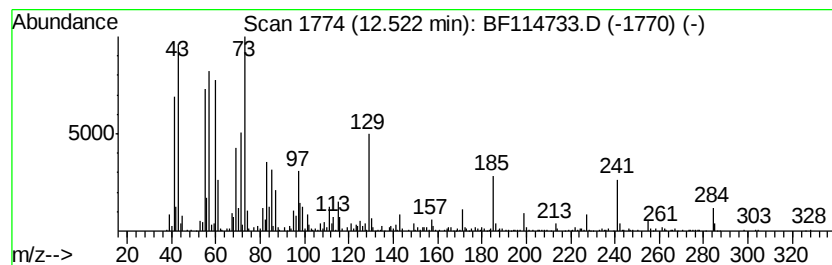
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Octadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.52	7.94 ng	817495	Chrysene-d12	13.82

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octadecanoic acid	284	C18H36O2	000057-11-4	99
2		Pentadecanoic acid	242	C15H30O2	001002-84-2	81
3		Tetradecanoic acid	228	C14H28O2	000544-63-8	74
4		n-Decanoic acid	172	C10H20O2	000334-48-5	70
5		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	64



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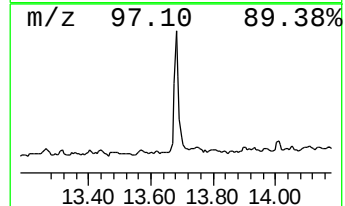
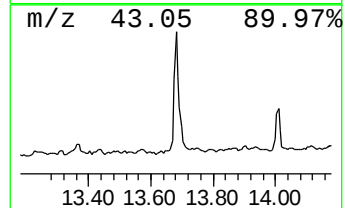
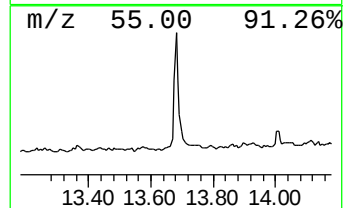
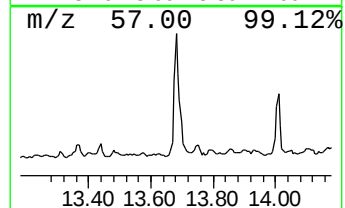
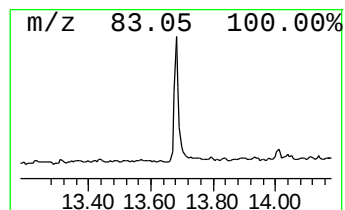
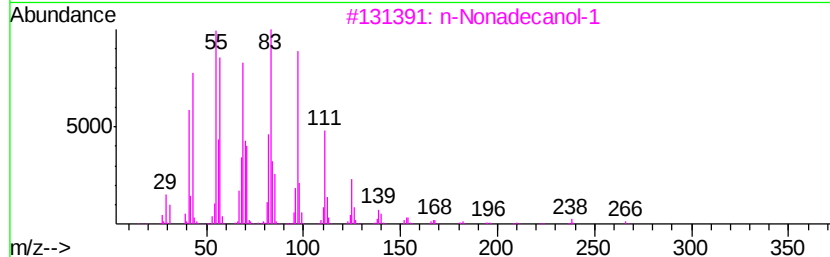
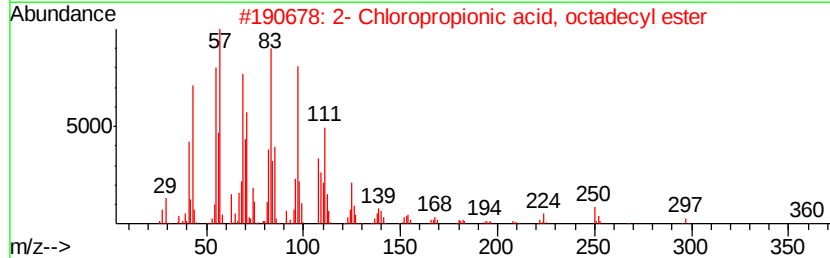
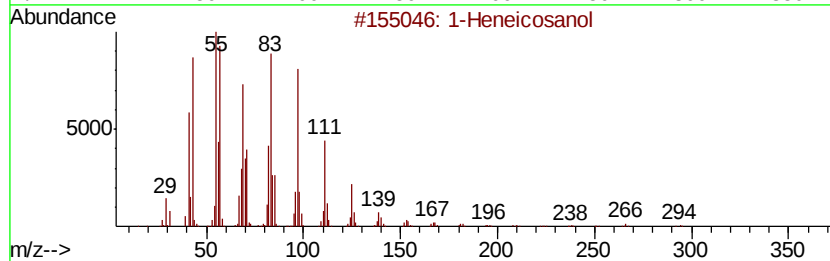
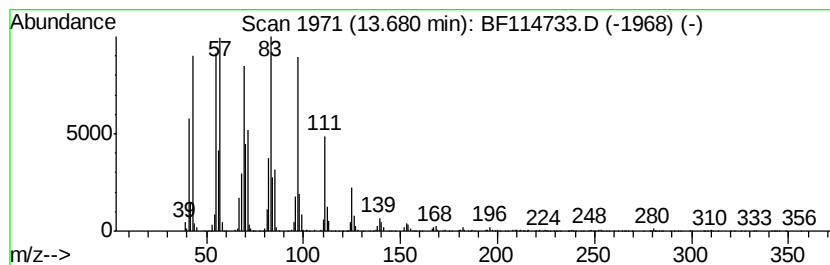
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 5 1-Heneicosanol Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.68	13.60 ng	1400180	Chrysene-d12	13.82	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heneicosanol	312	C21H44O	015594-90-8	95
2	2- Chloropropionic acid, octadec...	360	C21H41ClO2	088104-31-8	94
3	n-Nonadecanol-1	284	C19H40O	001454-84-8	94
4	Trifluoroacetic acid, pentadecyl...	324	C17H31F3O2	959010-23-2	93
5	Trifluoroacetoxyl hexadecane	338	C18H33F3O2	006222-03-3	93



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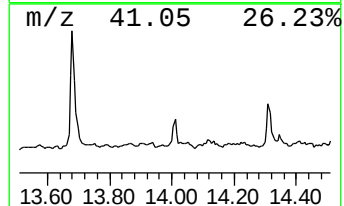
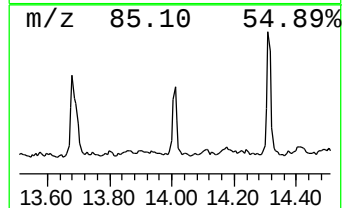
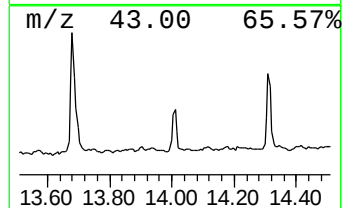
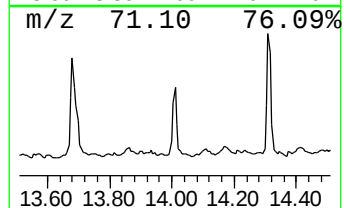
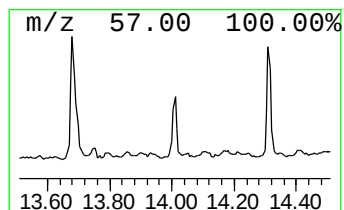
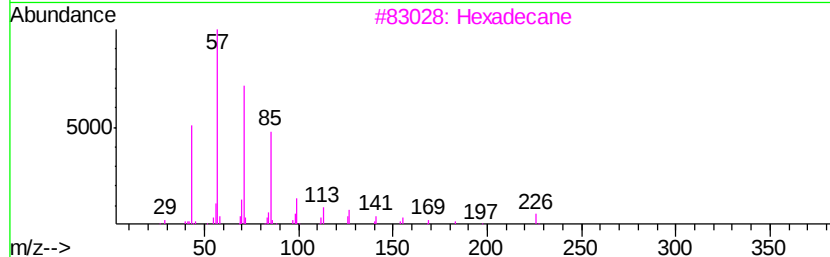
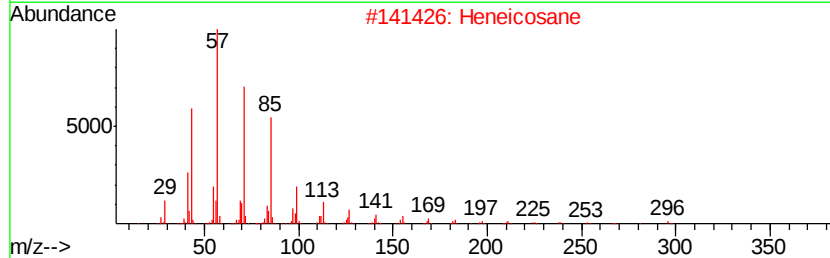
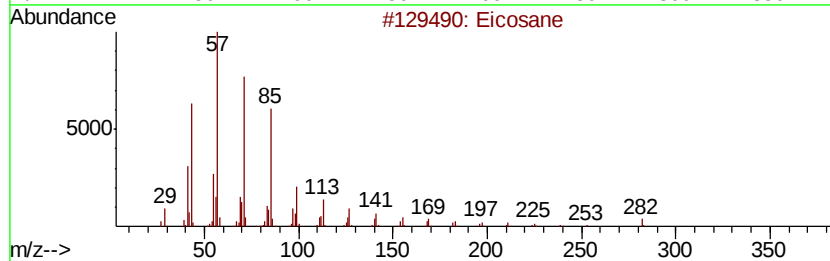
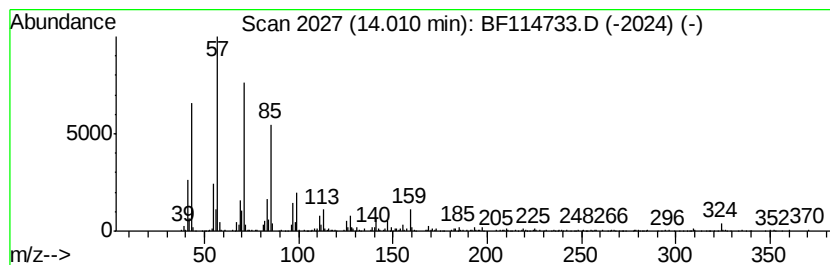
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Eicosane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.01	3.35 ng	344984	Chrysene-d12	13.82

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane		282	C20H42	000112-95-8	90
2	Heneicosane		296	C21H44	000629-94-7	87
3	Hexadecane		226	C16H34	000544-76-3	86
4	Tetratetracontane		619	C44H90	007098-22-8	83
5	Octacosane		394	C28H58	000630-02-4	76



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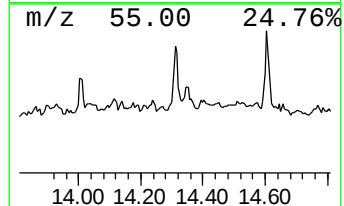
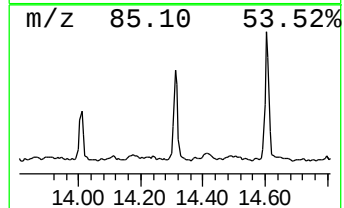
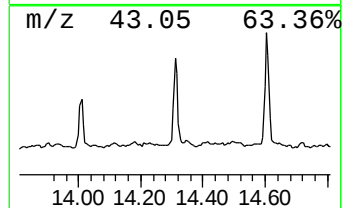
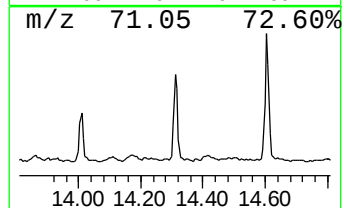
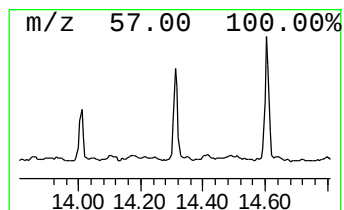
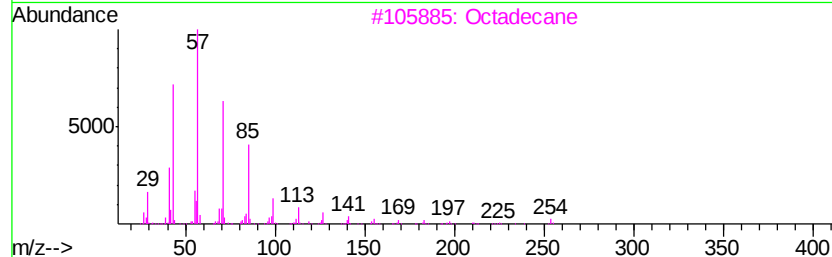
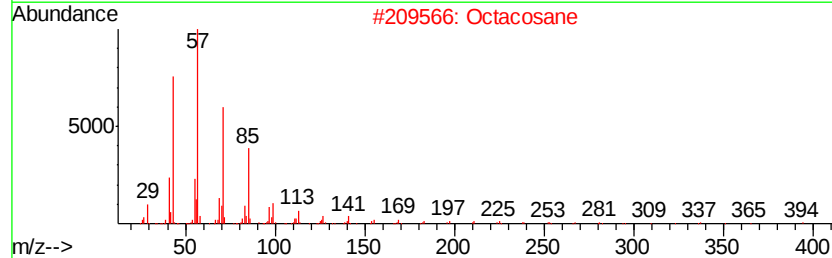
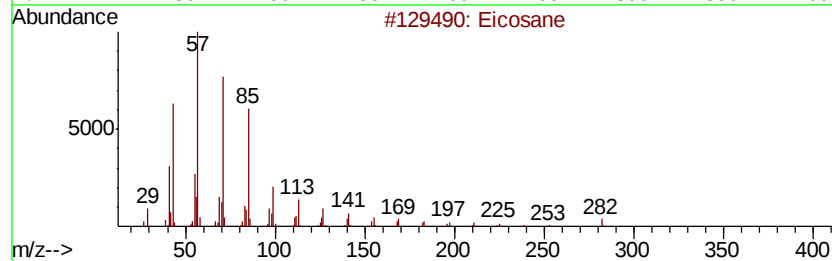
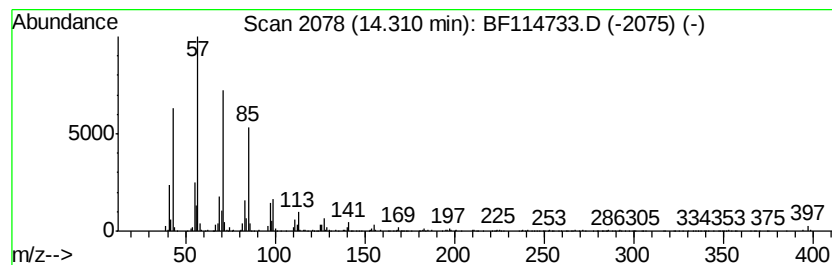
Instrument :
BNA_F
ClientSampleId :
TP-3

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF052919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Heneicosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD		R.T.	
14.31	6.11 ng	629381	Chrysene-d12		13.82	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane		282	C20H42	000112-95-8	97
2	Octacosane		394	C28H58	000630-02-4	93
3	Octadecane		254	C18H38	000593-45-3	93
4	Heneicosane		296	C21H44	000629-94-7	91
5	Hexatriacontane		507	C36H74	000630-06-8	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF053119\
Data File : BF114733.D
Acq On : 1 Jun 2019 3:13
Operator : HP/JU
Sample : K3072-03
Misc :
ALS Vial : 18 Sample Multiplier: 1

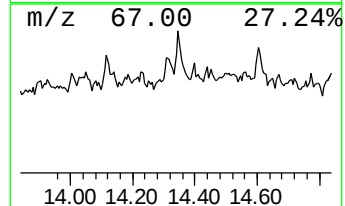
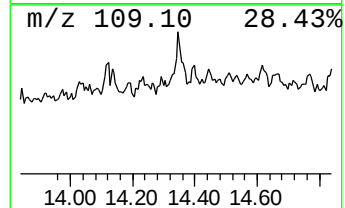
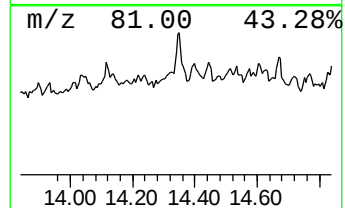
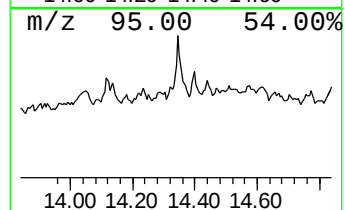
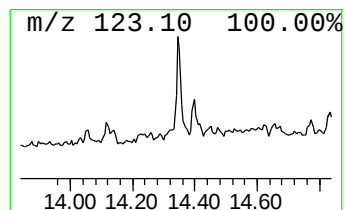
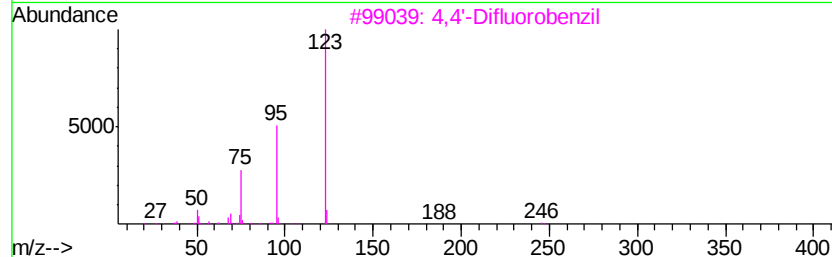
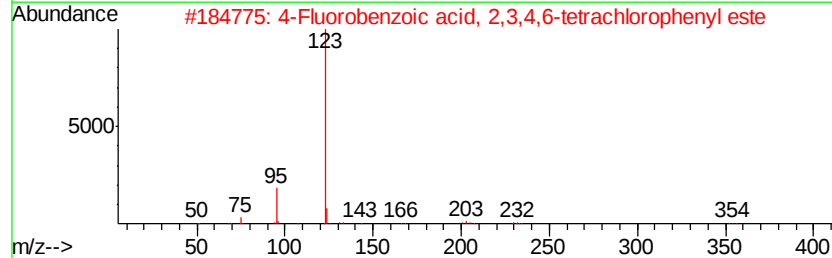
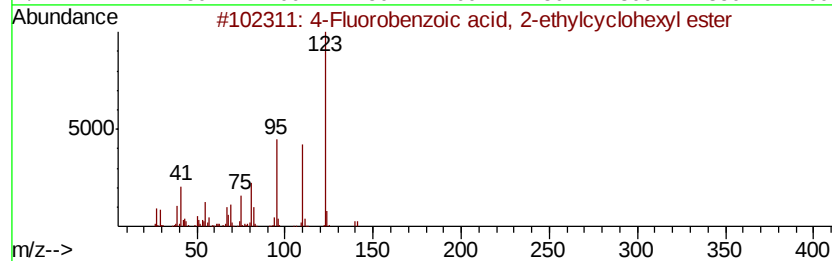
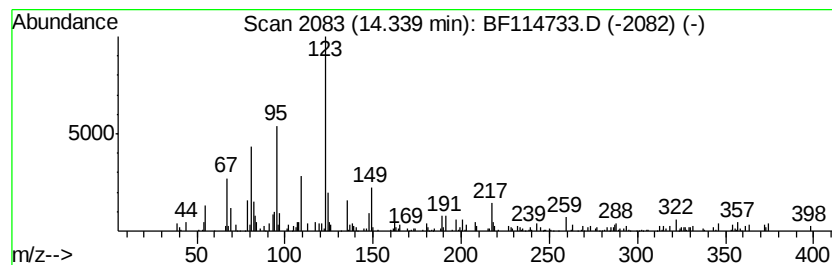
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF052919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 4-Fluorobenzoic acid, 2-eth... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.		
14.34	3.02 ng	310740	Chrysene-d12	13.82		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Fluorobenzoic acid, 2-ethylcvc...	250	C15H19F02	1000279-06-0	50	
2	4-Fluorobenzoic acid, 2,3,4,6-te...	352	C13H5Cl4F02	1000355-68-1	50	
3	4,4'-Difluorobenzil	246	C14H8F2O2	000579-39-5	49	
4	4-Fluorobenzoic acid, 3-methylbu...	208	C12H13F02	1000299-15-1	49	
5	Methyl 2-fluorobenzoate	154	C8H7F02	000394-35-4	47	



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF053119\
 Data File : BF114733.D
 Acq On : 1 Jun 2019 3:13
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 Sample : K3072-03
 Misc :
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Instrument :
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 ClientSampleId :
 TP-3

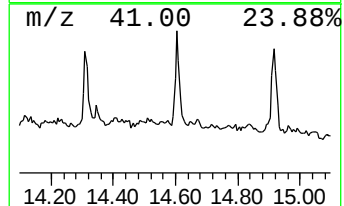
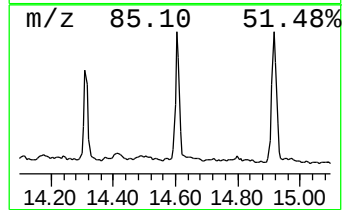
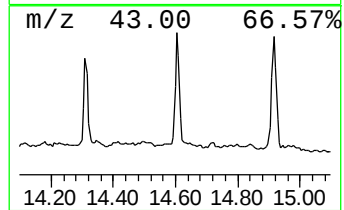
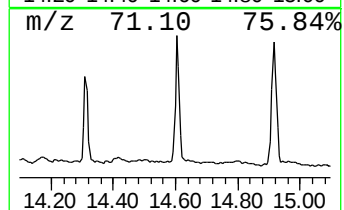
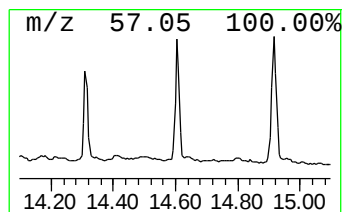
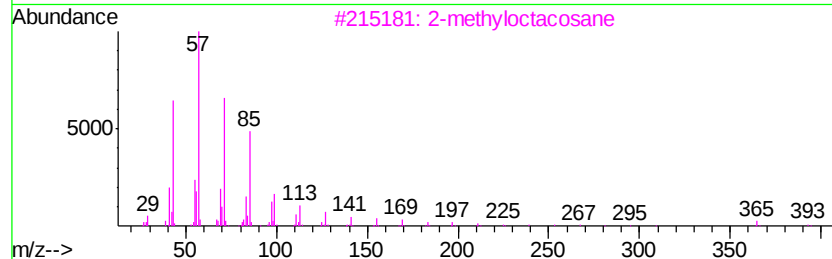
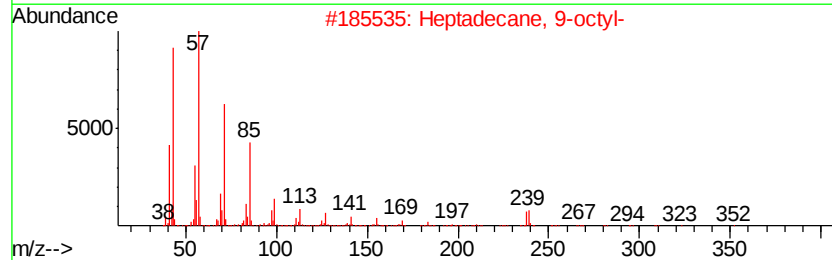
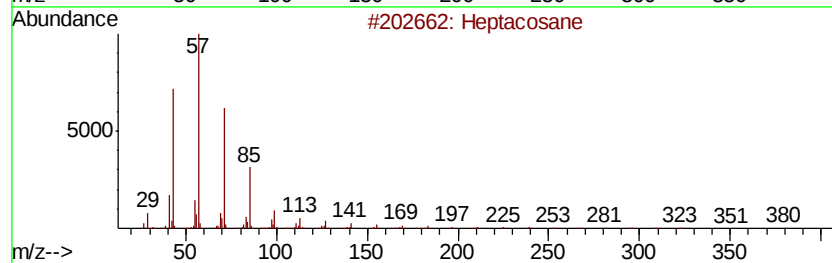
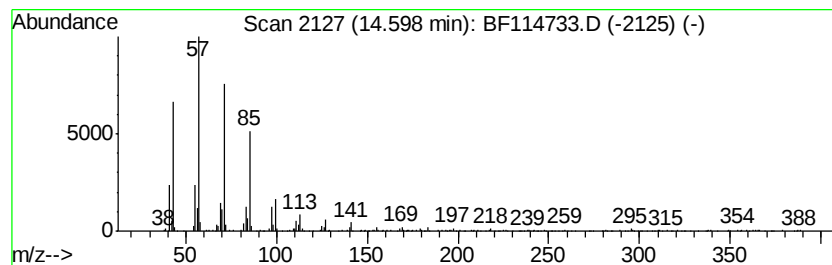
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 9 Heptacosane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.60	8.63 ng	861411	Perylene-d12	15.20

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptacosane		380	C27H56	000593-49-7	93
2	Heptadecane, 9-octyl-		352	C25H52	007225-64-1	91
3	2-methyloctacosane		408	C29H60	1000376-72-8	91
4	Heneicosane		296	C21H44	000629-94-7	91
5	Octacosane		394	C28H58	000630-02-4	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF053119\
Data File : BF114733.D
Acq On : 1 Jun 2019 3:13
Operator : HP/JU
Sample : K3072-03
Misc :
ALS Vial : 18 Sample Multiplier: 1

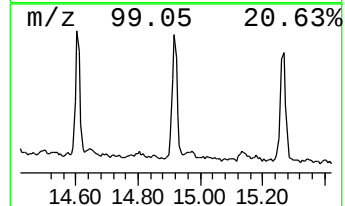
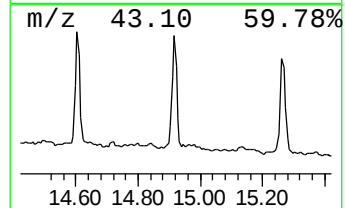
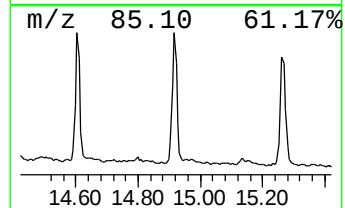
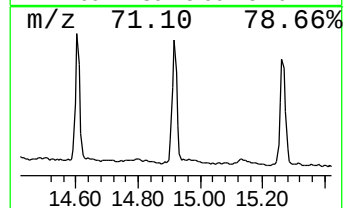
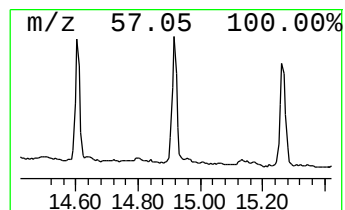
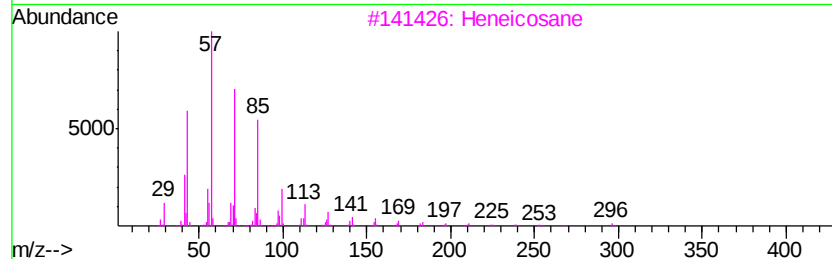
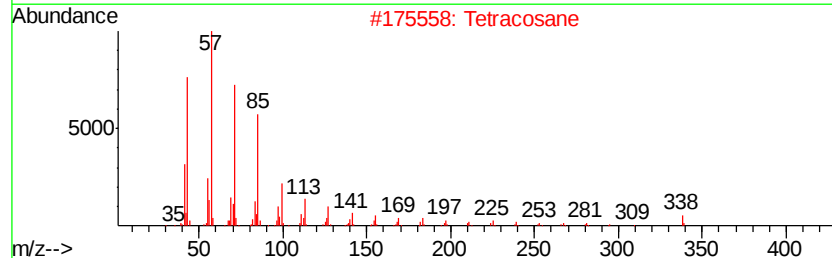
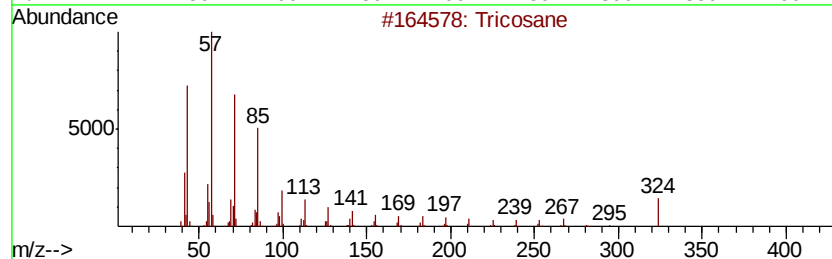
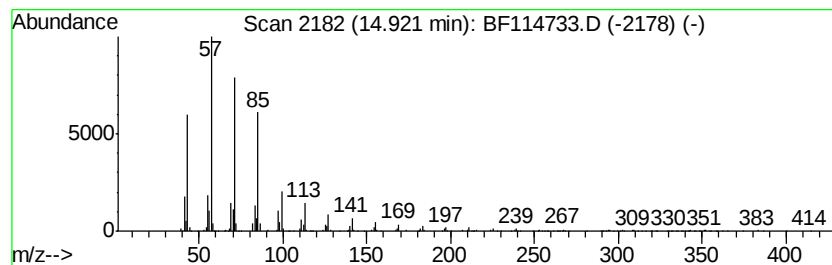
Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF052919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Tricosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.92	8.62 ng	859836	Perylene-d12	15.20
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Tricosane	324 C23H48	000638-67-5	95
2	Tetracosane	338 C24H50	000646-31-1	93
3	Heneicosane	296 C21H44	000629-94-7	91
4	Heptadecane, 9-hexyl-	324 C23H48	055124-79-3	91
5	Eicosane, 7-hexyl-	366 C26H54	055333-99-8	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF053119\
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Sample : K3072-03
Misc :
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Instrument :
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ClientSampleId :
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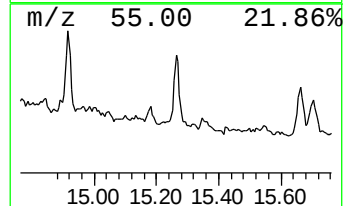
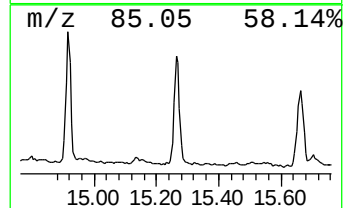
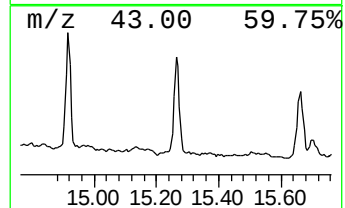
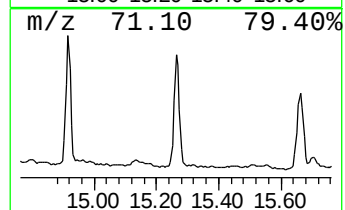
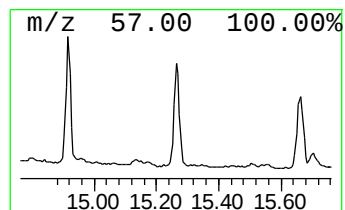
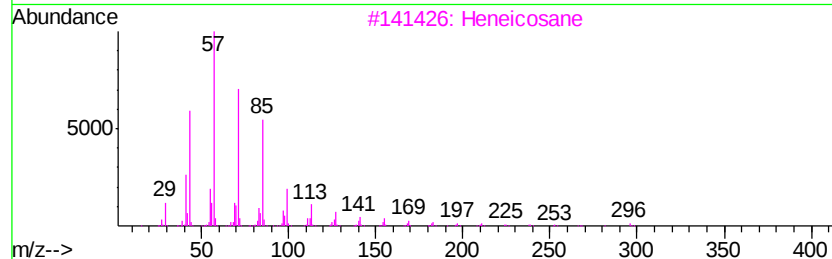
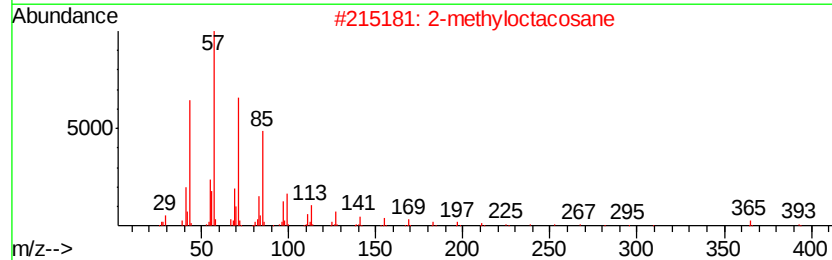
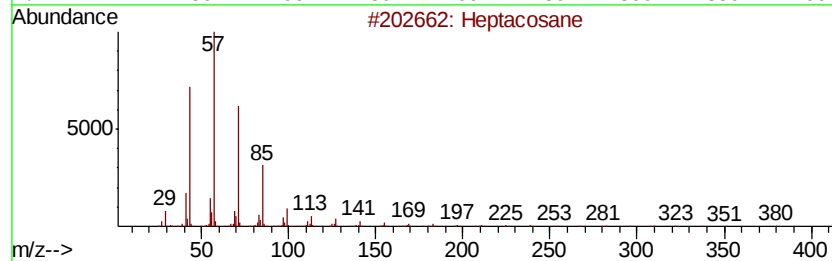
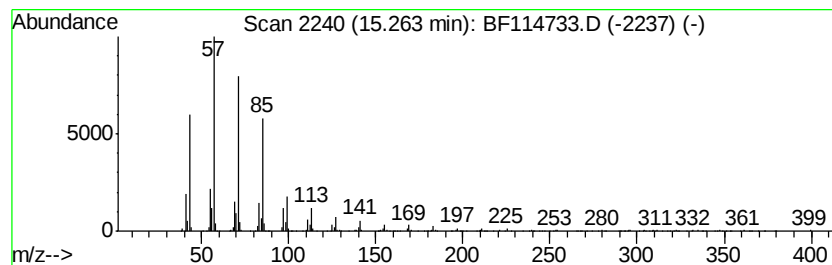
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Tridecane, 1-iodo- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.26	9.18 ng	915523	Perylene-d12	15.20

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptacosane	380	C27H56	000593-49-7	93
2		2-methyloctacosane	408	C29H60	1000376-72-8	91
3		Heneicosane	296	C21H44	000629-94-7	91
4		Heptadecane	240	C17H36	000629-78-7	91
5		Tridecane, 1-iodo-	310	C13H27I	035599-77-0	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF053119\
Data File : BF114733.D
Acq On : 1 Jun 2019 3:13
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Sample : K3072-03
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TP-3

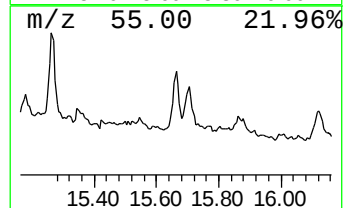
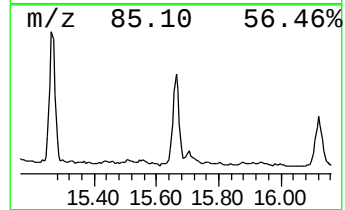
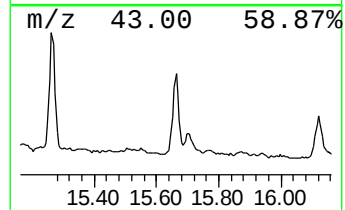
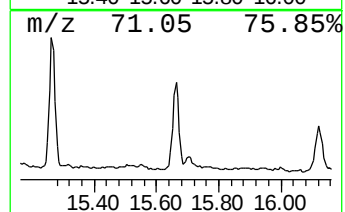
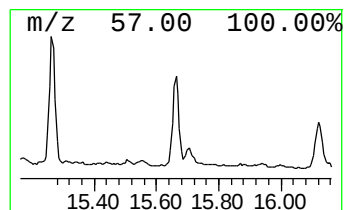
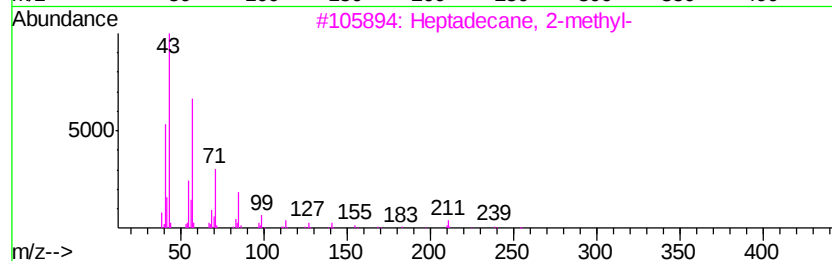
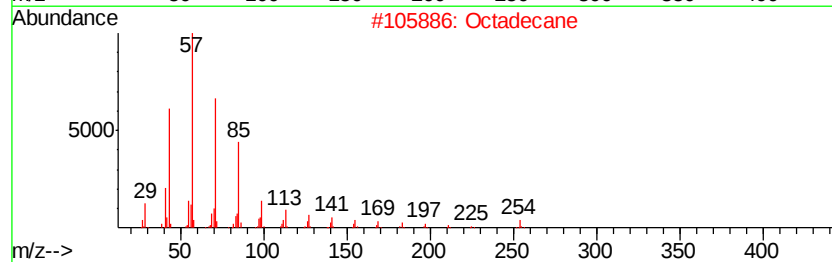
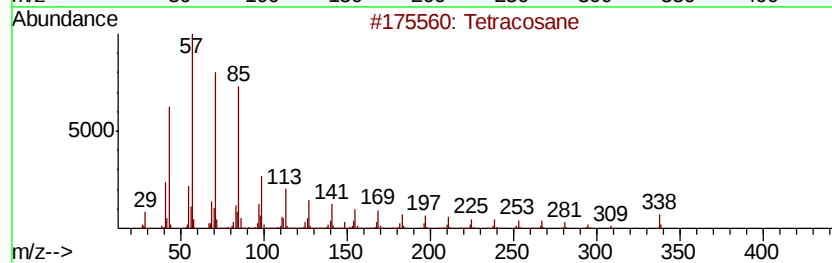
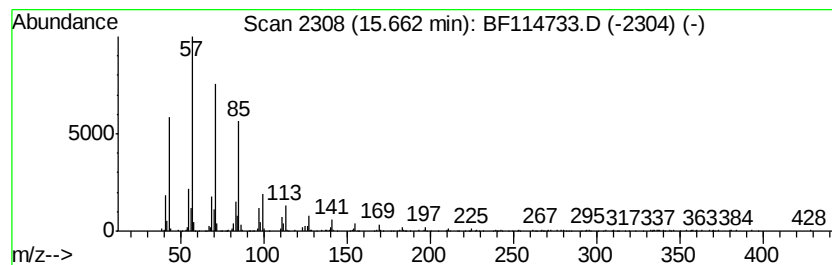
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Heptadecane, 2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.66	6.56 ng	654133	Perylene-d12	15.20

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetracosane	338	C24H50	000646-31-1	95
2			Octadecane	254	C18H38	000593-45-3	94
3			Heptadecane, 2-methyl-	254	C18H38	001560-89-0	91
4			Heneicosane	296	C21H44	000629-94-7	91
5			2-methyloctacosane	408	C29H60	1000376-72-8	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF053119\
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ALS Vial : 18 Sample Multiplier: 1

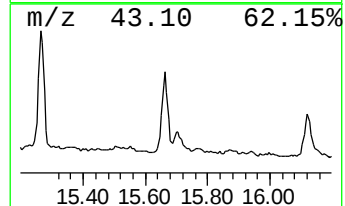
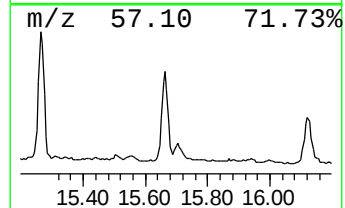
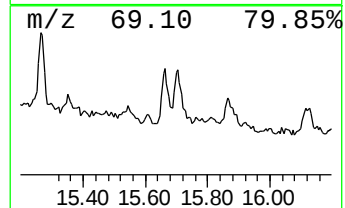
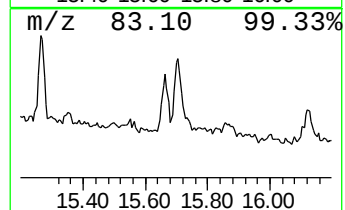
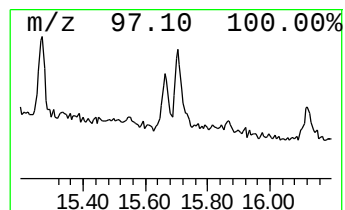
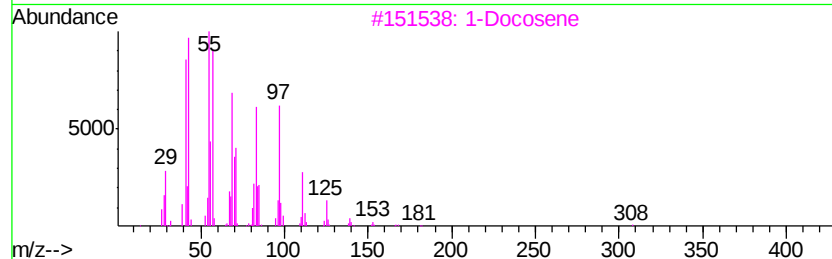
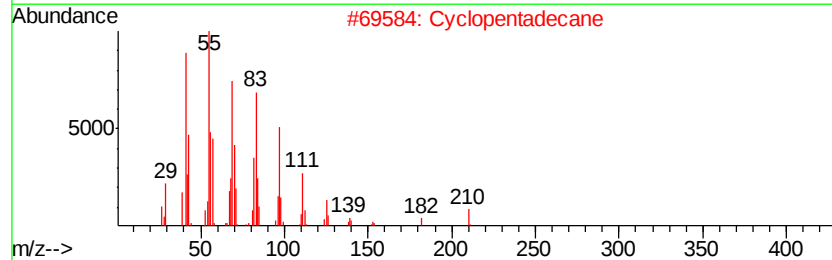
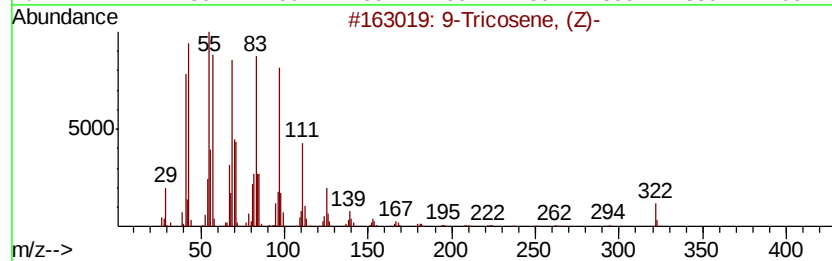
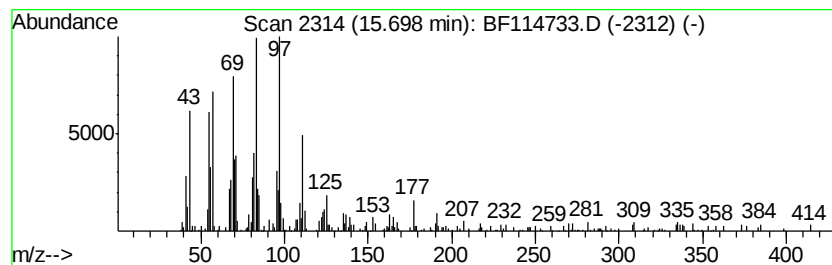
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 9-Tricosene, (Z)- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.70	4.19 ng	417872	Perylene-d12	15.20
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		9-Tricosene, (Z)-	322 C23H46	027519-02-4 93
2		Cyclopentadecane	210 C15H30	000295-48-7 87
3		1-Docosene	308 C22H44	001599-67-3 86
4		13-Tetradecen-1-ol acetate	254 C16H30O2	056221-91-1 83
5		Behenic alcohol	326 C22H46O	000661-19-8 81



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF053119\
Data File : BF114733.D
Acq On : 1 Jun 2019 3:13
Operator : HP/JU
Sample : K3072-03
Misc :
ALS Vial : 18 Sample Multiplier: 1

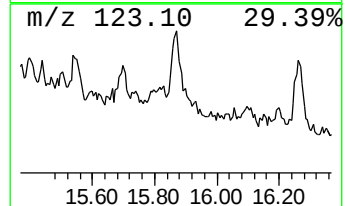
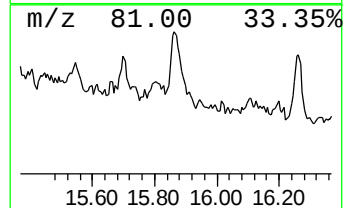
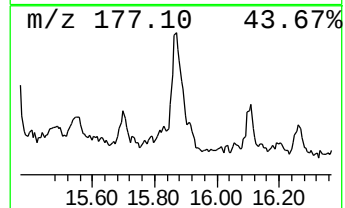
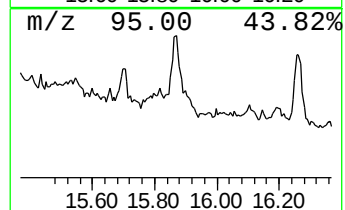
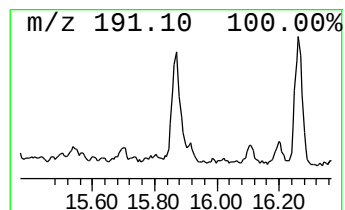
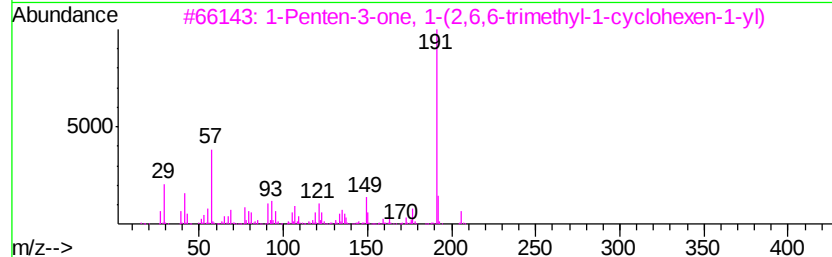
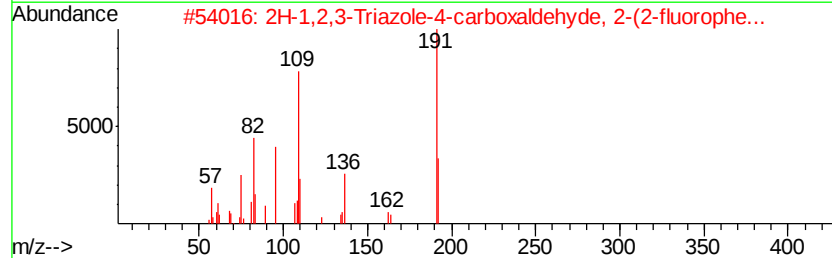
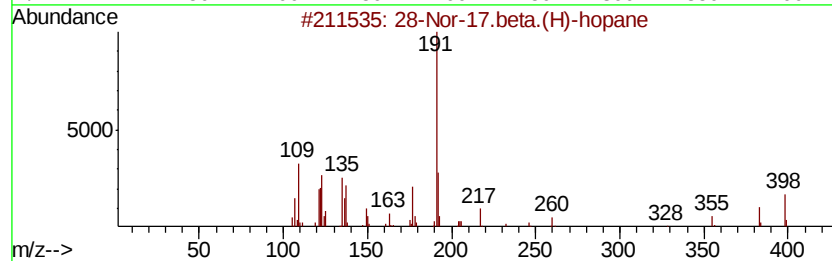
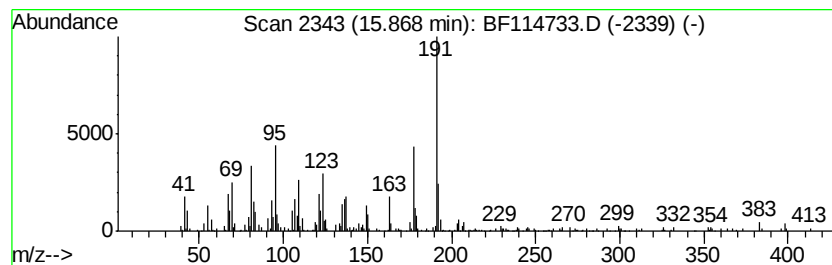
Instrument :
BNA_F
ClientSampleId :
TP-3

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF052919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 28-Nor-17.beta.(H)-hopane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.87	5.61 ng	560064	Perylene-d12	15.20
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	28-Nor-17.beta.(H)-hopane	398 C29H50	036728-72-0	60
2	2H-1,2,3-Triazole-4-carboxaldehy...	191 C9H6FN3O	051306-43-5	43
3	1-Penten-3-one, 1-(2,6,6-trimeth...	206 C14H22O	000127-43-5	38
4	5-(p-Aminophenyl)-2-thiazolamine	191 C9H9N3S	090349-87-4	27
5	17.alfa.,21.beta.-28,30-Bisnorho...	384 C28H48	1000360-26-1	27



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF053119\
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Sample : K3072-03
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TP-3

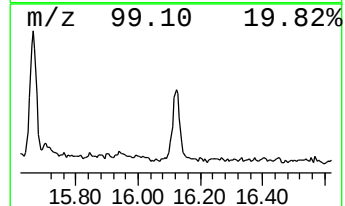
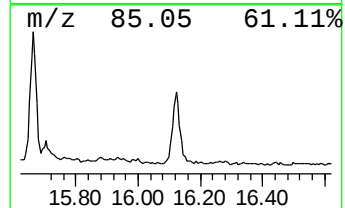
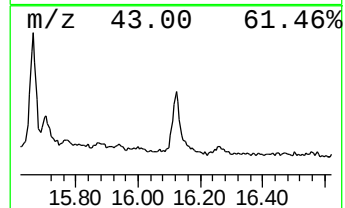
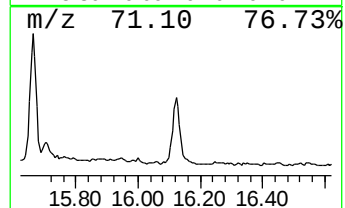
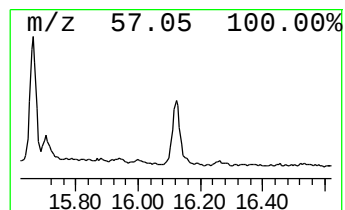
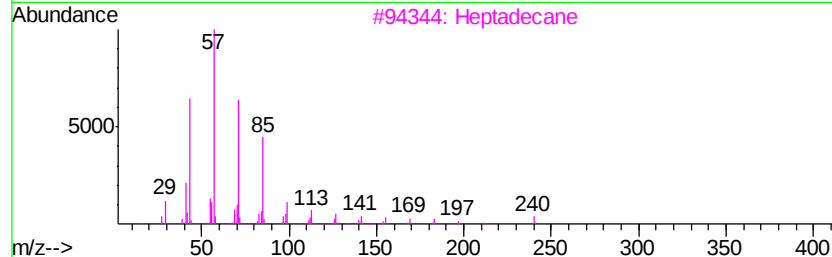
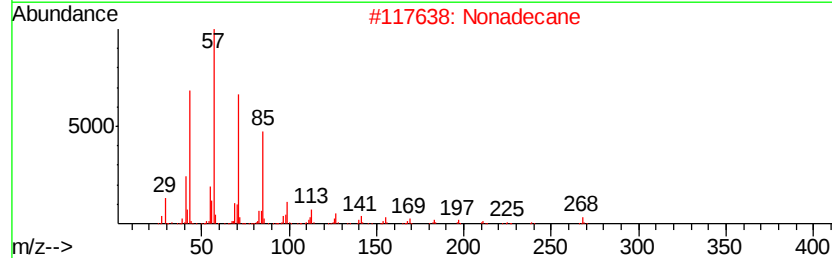
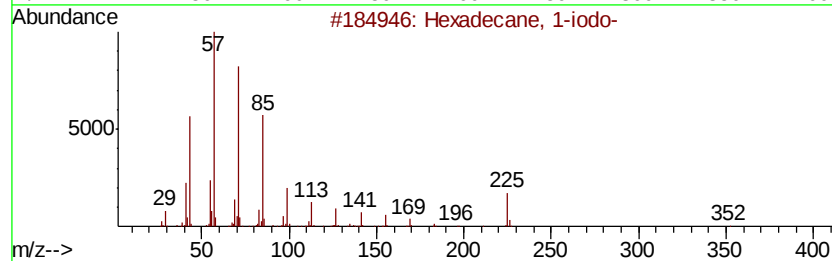
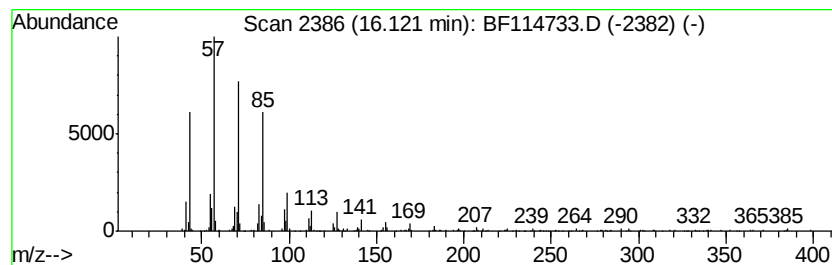
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF052919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Hexadecane, 1-iodo- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.12	3.13 ng	312469	Perylene-d12	15.20

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	96
2		Nonadecane	268	C19H40	000629-92-5	94
3		Heptadecane	240	C17H36	000629-78-7	91
4		Heneicosane	296	C21H44	000629-94-7	91
5		Hentriacontane	437	C31H64	000630-04-6	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF053119\
Data File : BF114733.D
Acq On : 1 Jun 2019 3:13
Operator : HP/JU
Sample : K3072-03
Misc :
ALS Vial : 18 Sample Multiplier: 1

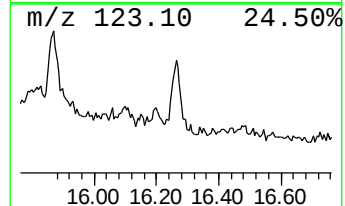
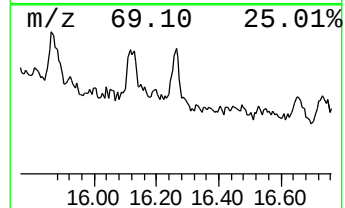
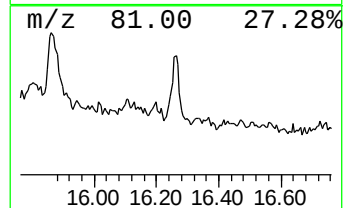
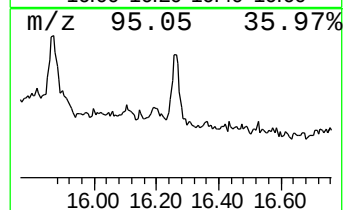
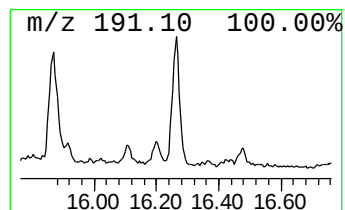
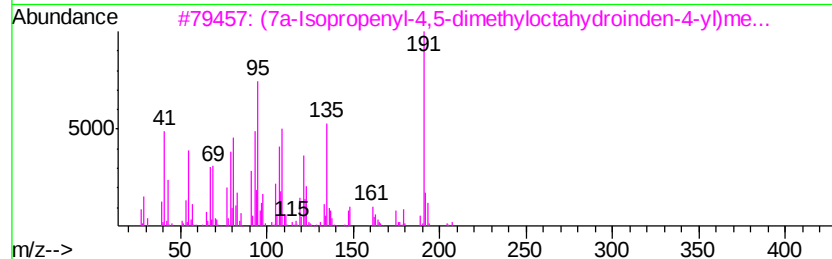
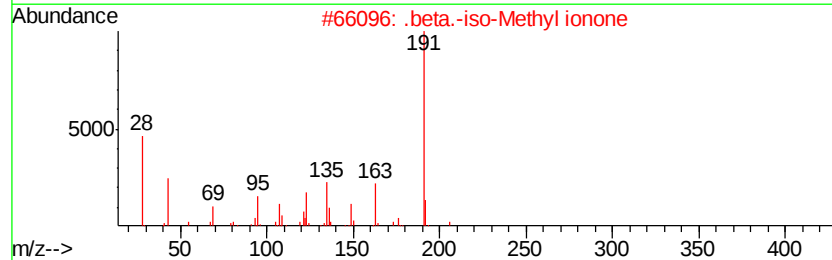
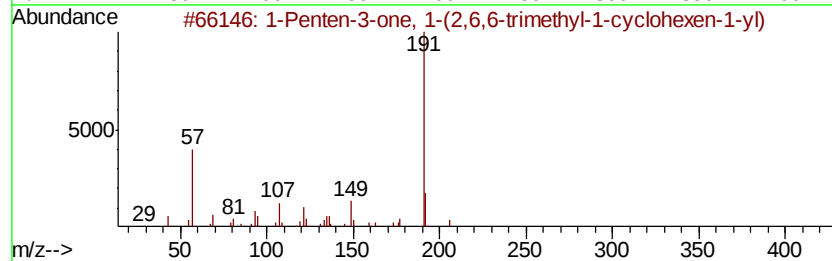
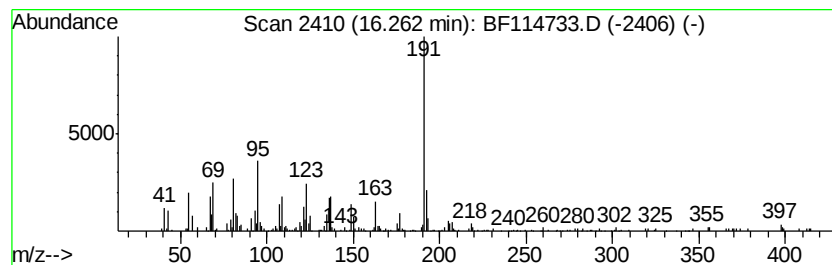
Instrument :
BNA_F
ClientSampleId :
TP-3

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF052919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 1-Penten-3-one, 1-(2,6,6-tr... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.26	3.71 ng	369949	Perylene-d12	15.20
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1-Penten-3-one, 1-(2,6,6-trimeth...	206 C14H22O	000127-43-5	70
2	.beta.-iso-Methyl ionone	206 C14H22O	1000285-40-2	68
3	(7a-Isopropenyl-4,5-dimethyl-octa...	222 C15H26O	1000187-02-9	50
4	5-(p-Aminophenyl)-2-thiazolamine	191 C9H9N3S	090349-87-4	47
5	Phenol, 3,5-bis(1,1-dimethylethyl)-	206 C14H22O	001138-52-9	38



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF053119\
 Data File : BF114733.D
 Acq On : 1 Jun 2019 3:13
 Operator : HP/JU
 Sample : K3072-03
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-3

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF052919.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
unknown6.47	6.47	83.8	ng	9850560	1	6.70	2351360	20.0
n-Hexadecanoic acid	11.74	12.0	ng	1532950	4	11.19	2564510	20.0
Octadecanoic acid	12.52	7.9	ng	817495	5	13.82	2058570	20.0
1-Heneicosanol	13.68	13.6	ng	1400180	5	13.82	2058570	20.0
Eicosane	14.01	3.4	ng	344984	5	13.82	2058570	20.0
Heneicosane	14.31	6.1	ng	629381	5	13.82	2058570	20.0
4-Fluorobenzoic a...	14.34	3.0	ng	310740	5	13.82	2058570	20.0
Heptacosane	14.60	8.6	ng	861411	6	15.20	1995400	20.0
Tricosane	14.92	8.6	ng	859836	6	15.20	1995400	20.0
Tridecane, 1-iodo-	15.26	9.2	ng	915523	6	15.20	1995400	20.0
Heptadecane, 2-me...	15.66	6.6	ng	654133	6	15.20	1995400	20.0
9-Tricosene, (Z)-	15.70	4.2	ng	417872	6	15.20	1995400	20.0
28-Nor-17.beta.(H...	15.87	5.6	ng	560064	6	15.20	1995400	20.0
Hexadecane, 1-iodo-	16.12	3.1	ng	312469	6	15.20	1995400	20.0
1-Penten-3-one, 1...	16.26	3.7	ng	369949	6	15.20	1995400	20.0