

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM031419\
 Data File : BM019169.D
 Acq On : 14 Mar 2019 15:03
 Operator : JU/SJ
 Sample : K1866-01
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 MR-MH-04-COMP

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_M\METHODS\8270-BM031119.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	5.258	379	386	404	rBV	4317971	6423288	37.03%	6.810%
2	6.840	647	655	674	rBV	4780064	7840819	45.20%	8.313%
3	7.193	708	715	735	rBV	4764674	7494455	43.20%	7.946%
4	7.657	787	794	803	rBV	738439	1194834	6.89%	1.267%
5	7.969	840	847	861	rBV	3124372	4981980	28.72%	5.282%
6	8.828	985	993	1012	rBV	2485764	4398423	25.36%	4.663%
7	10.440	1260	1267	1281	rBV	954808	1710349	9.86%	1.813%
8	12.922	1682	1689	1711	rBV	7322181	10807213	62.30%	11.458%
9	13.186	1729	1734	1742	rVB	181332	296958	1.71%	0.315%
10	13.775	1829	1834	1845	rBV	377885	605521	3.49%	0.642%
11	14.310	1918	1925	1944	rBV2	1709427	2645103	15.25%	2.804%
12	15.810	2173	2180	2199	rBV	6983464	10011089	57.71%	10.614%
13	17.063	2387	2393	2406	rBV	2297354	3414570	19.68%	3.620%
14	17.957	2539	2545	2561	rBV	830783	1375122	7.93%	1.458%
15	18.816	2686	2691	2698	rBV	938906	1314144	7.58%	1.393%
16	19.139	2741	2746	2750	rVB2	197462	265146	1.53%	0.281%
17	19.245	2759	2764	2771	rBV	260635	449137	2.59%	0.476%
18	19.704	2836	2842	2850	rBV	14456802	17346323	100.00%	18.390%
19	19.998	2889	2892	2897	rVB	214333	248821	1.43%	0.264%
20	20.498	2973	2977	2981	rBV2	237742	268397	1.55%	0.285%
21	20.962	3052	3056	3068	rVB	1051325	1354380	7.81%	1.436%
22	21.262	3102	3107	3119	rVB	3022032	3855464	22.23%	4.088%
23	21.398	3126	3130	3137	rBV	353237	403435	2.33%	0.428%
24	21.827	3199	3203	3210	rBV	423616	536048	3.09%	0.568%
25	22.286	3277	3281	3288	rBV2	427412	591640	3.41%	0.627%
26	22.786	3362	3366	3371	rBV	337420	465200	2.68%	0.493%
27	23.351	3458	3462	3467	rVB	279417	443454	2.56%	0.470%
28	23.498	3480	3487	3501	rBV	1712727	3228004	18.61%	3.422%
29	23.992	3567	3571	3579	rVB	201867	352961	2.03%	0.374%

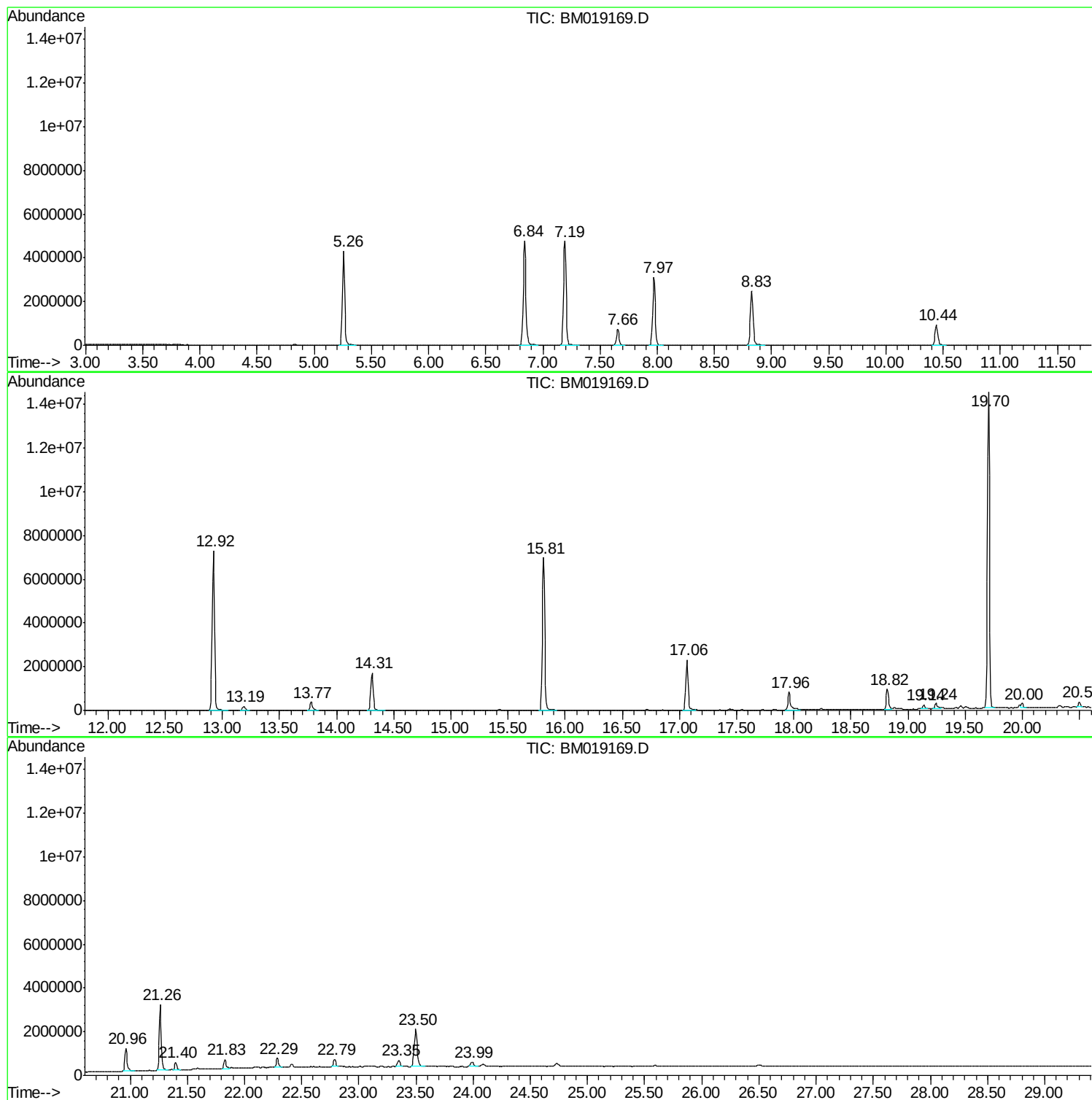
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM031119.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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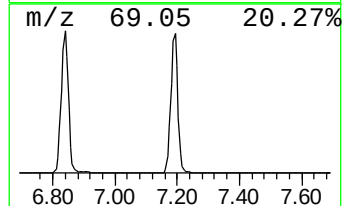
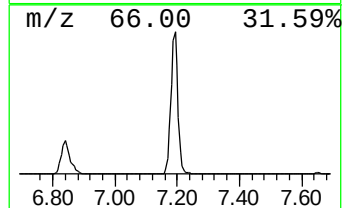
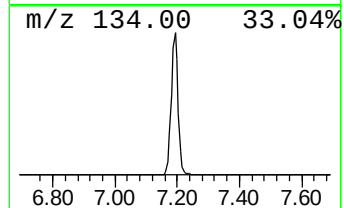
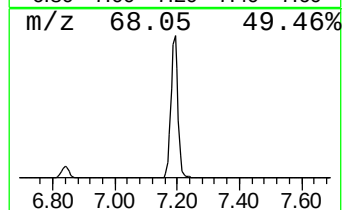
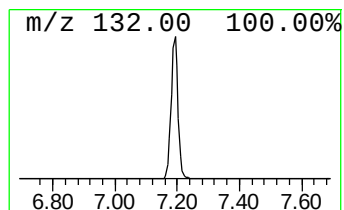
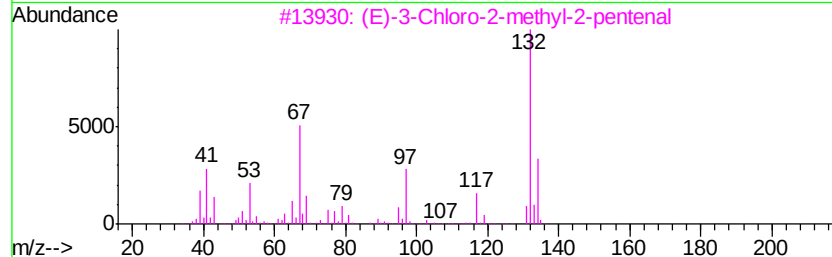
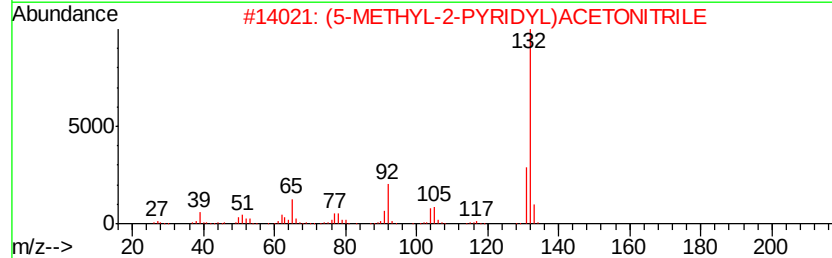
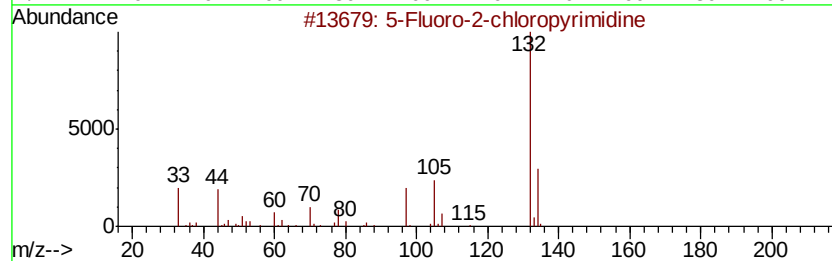
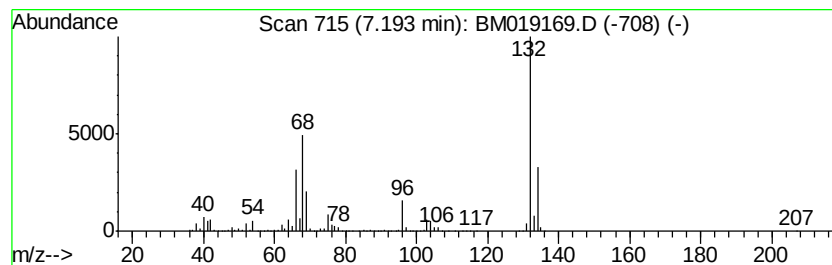
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 unknown7.19 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.19	125.45 ng	7494460	1,4-Dichlorobenzene-d4	7.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
2		(5-METHYL-2-PYRIDYL)ACETONITRILE	132	C8H8N2	1000241-93-9	10
3		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9



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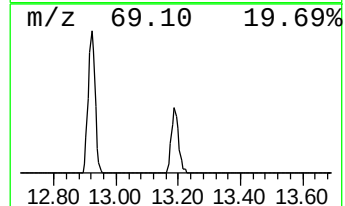
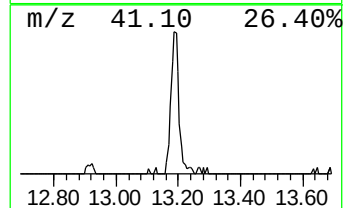
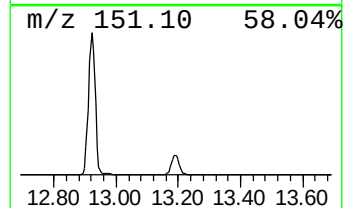
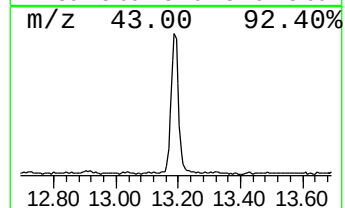
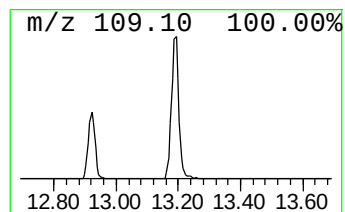
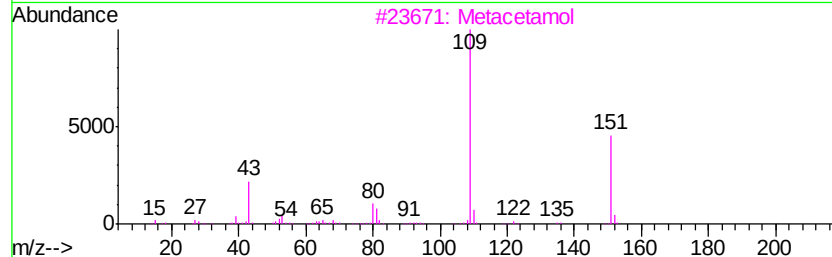
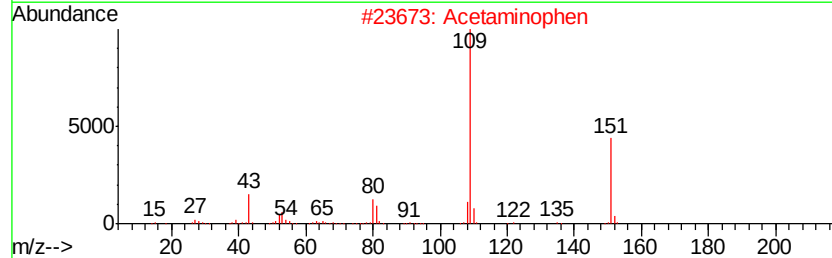
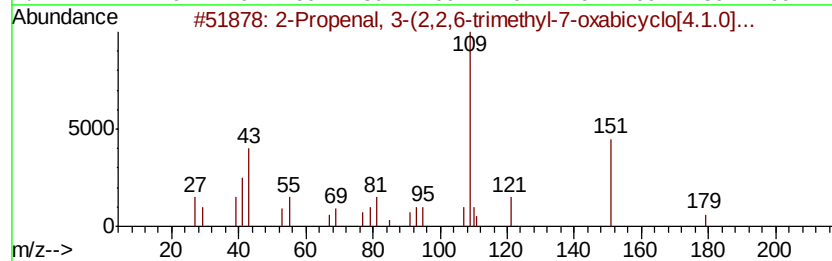
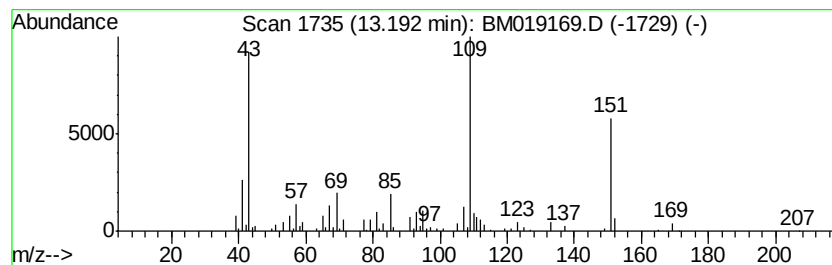
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Peak Number 3 2-Propenal, 3-(2,2,6-trimet... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD		R.T.
13.19	2.25 ng	296958	Acenaphthene-d10		14.31
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propenal, 3-(2,2,6-trimethyl-7...	194	C12H18O2	055759-91-6	64
2	Acetaminophen	151	C8H9NO2	000103-90-2	58
3	Metacetamol	151	C8H9NO2	000621-42-1	58
4	m-Isopropoxvaniline	151	C9H13NO	041406-00-2	53
5	2,4-Heptadienal, 2,4-dimethyl-	138	C9H14O	042452-48-2	43



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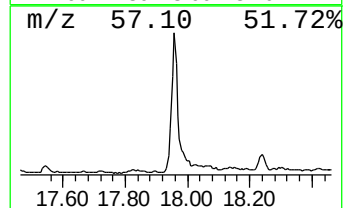
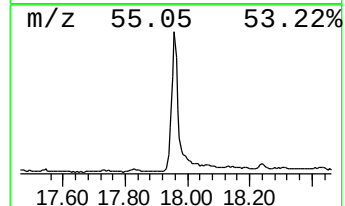
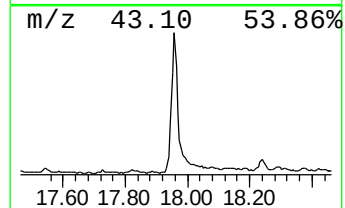
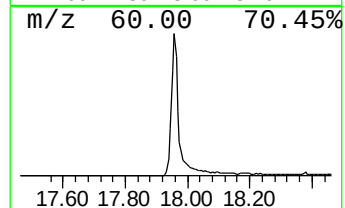
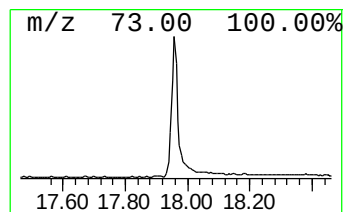
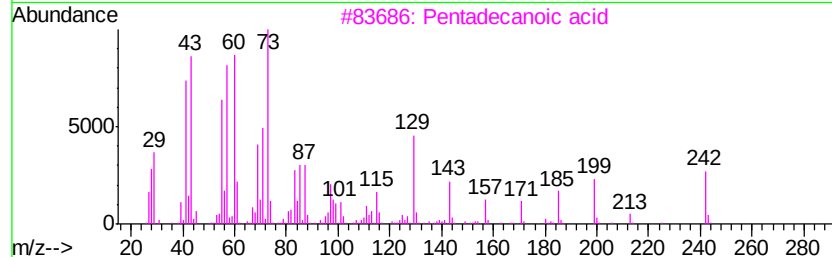
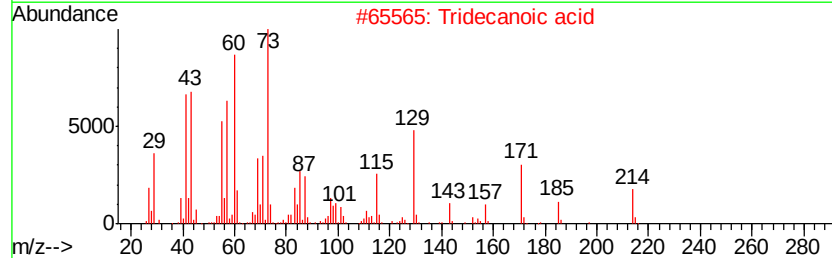
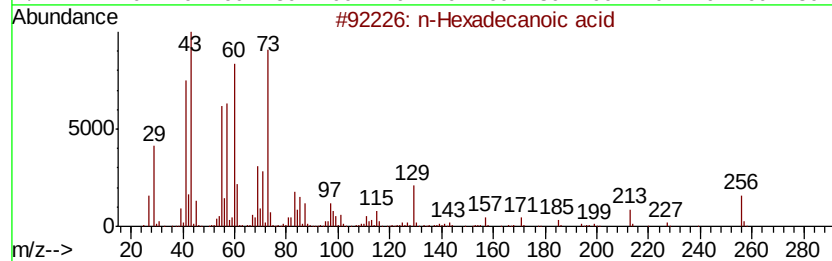
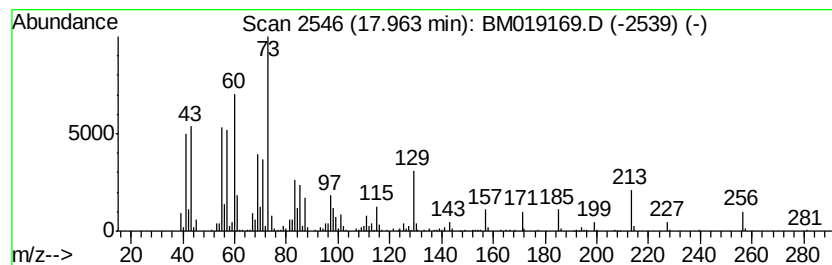
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Peak Number 4 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.	
17.96	8.05 ng	1375120	Phenanthrene-d10	17.06	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
2	Tridecanoic acid	214	C13H26O2	000638-53-9	91
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	72
4	Dodecane	170	C12H26	000112-40-3	18
5	Tetradecane, 1-iodo-	324	C14H29I	019218-94-1	18



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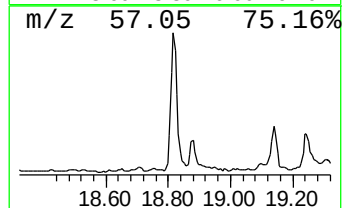
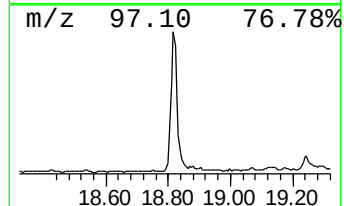
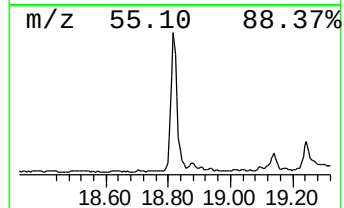
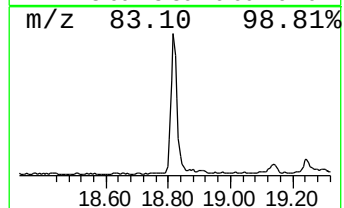
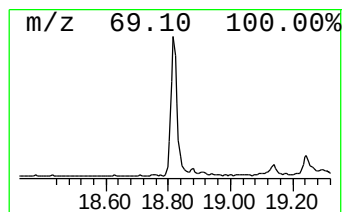
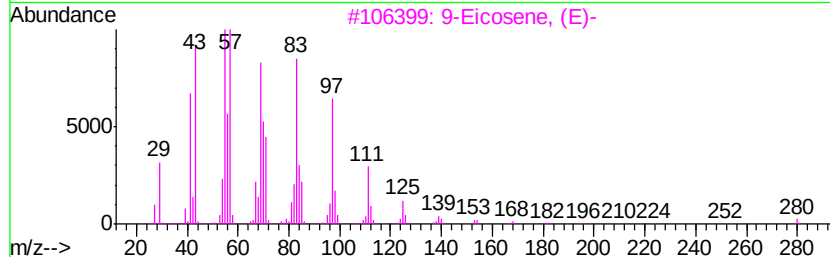
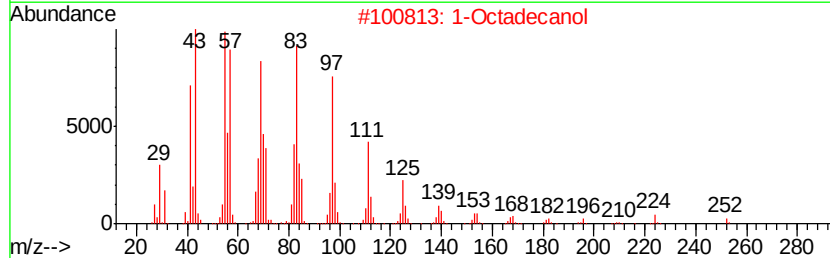
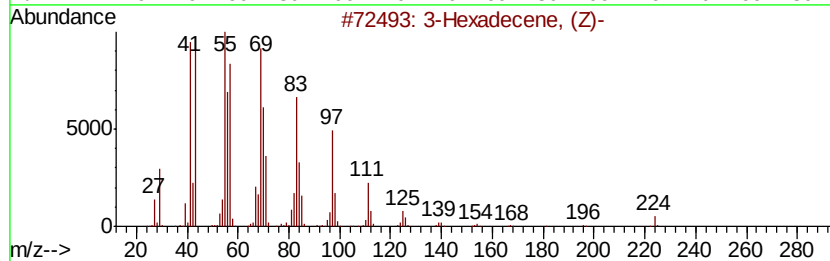
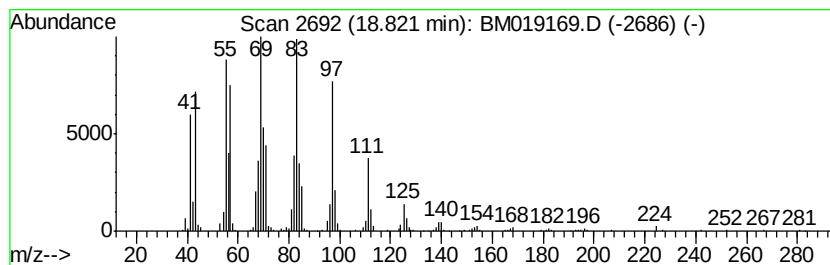
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Peak Number 5 3-Hexadecene, (Z)- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.82	7.70 ng	1314140	Phenanthrene-d10	17.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Hexadecene, (Z)-	224	C16H32	034303-81-6	95
2		1-Octadecanol	270	C18H38O	000112-92-5	91
3		9-Eicosene, (E)-	280	C20H40	074685-29-3	91
4		3-Eicosene, (E)-	280	C20H40	074685-33-9	91
5		1-Hexadecanol	242	C16H34O	036653-82-4	91



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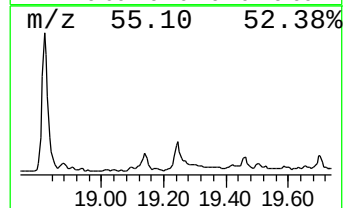
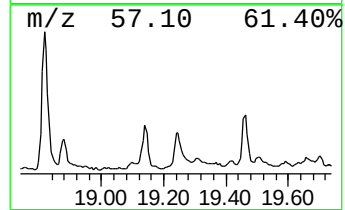
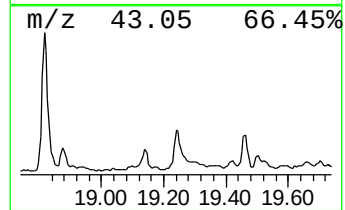
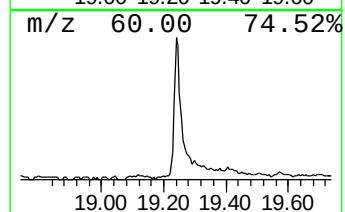
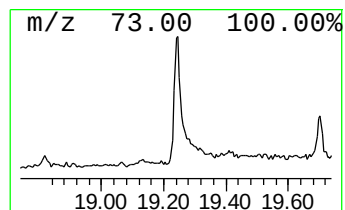
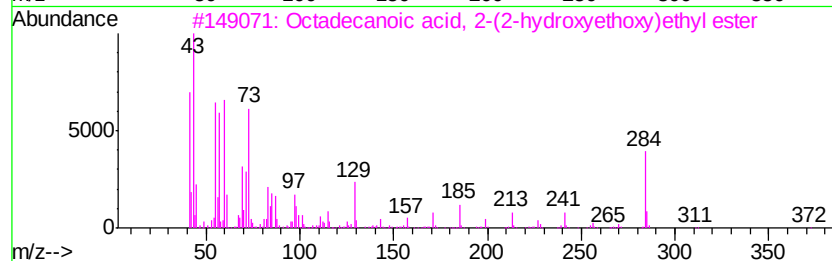
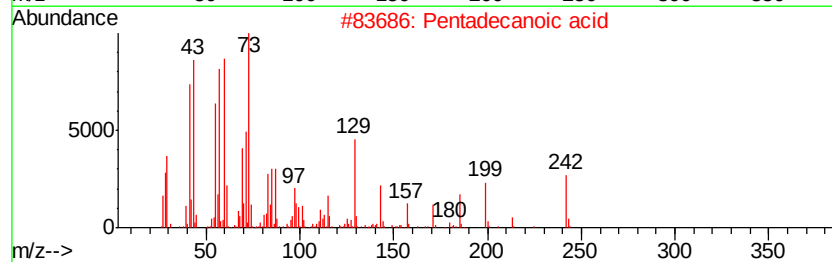
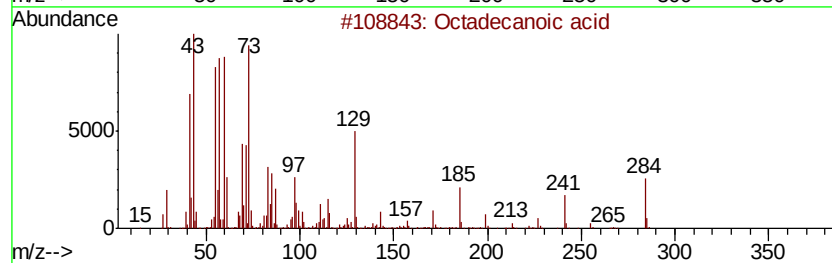
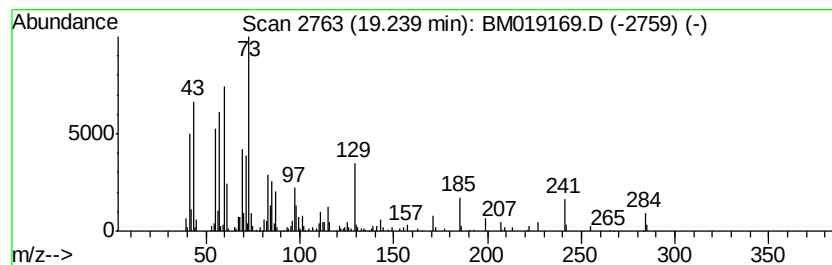
Instrument :
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Peak Number 6 Octadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.24	2.33 ng	449137	Chrysene-d12	21.26	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	96
2	Pentadecanoic acid	242	C15H30O2	001002-84-2	86
3	Octadecanoic acid, 2-(2-hydroxyve...	372	C22H44O4	000106-11-6	72
4	Tetradecanoic acid	228	C14H28O2	000544-63-8	64
5	n-Decanoic acid	172	C10H20O2	000334-48-5	38



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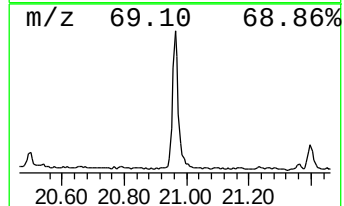
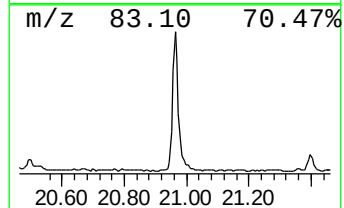
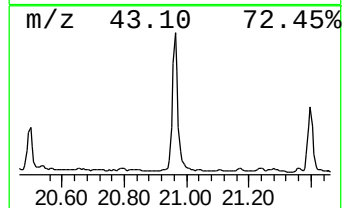
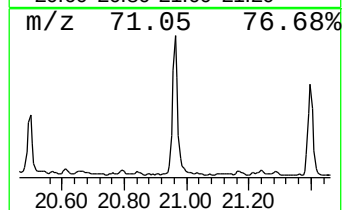
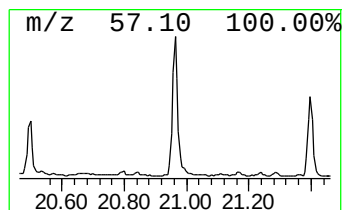
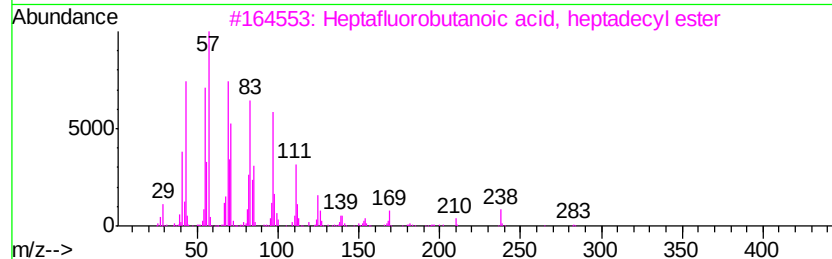
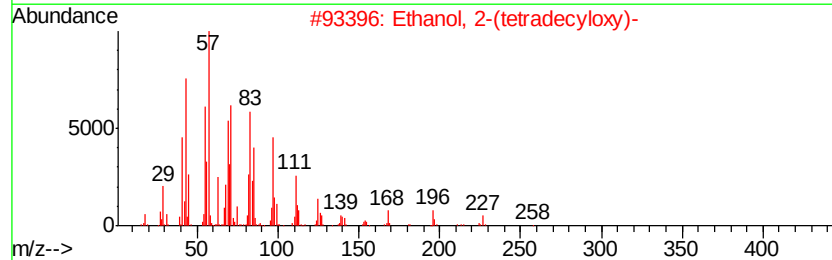
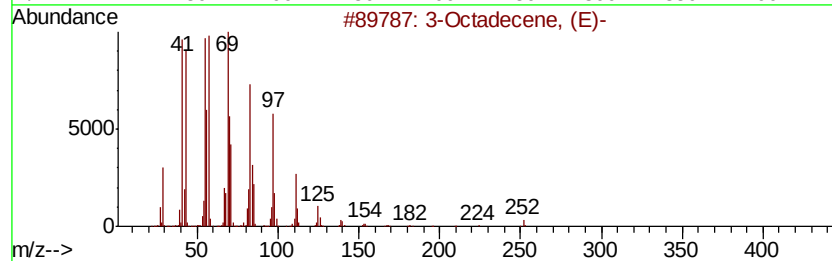
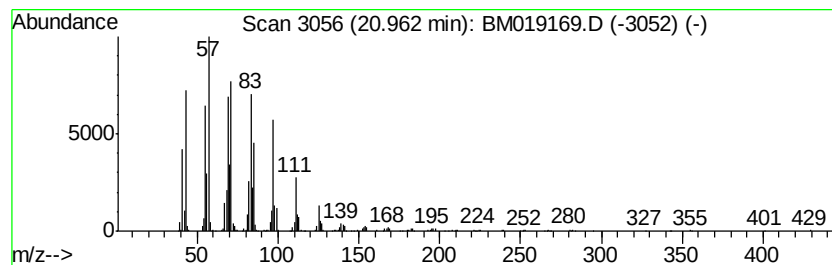
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 3-Octadecene, (E)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.96	7.03 ng	1354380	Chrysene-d12	21.26

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Octadecene, (E)-	252	C18H36	007206-19-1	93
2		Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	91
3		Heptafluorobutanoic acid, heptad...	452	C21H35F7O2	1000282-97-3	76
4		Pentafluoropropionic acid, hepta...	402	C20H35F5O2	1000283-04-2	76
5		Cyclotetradecane	196	C14H28	000295-17-0	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM031419\
Data File : BM019169.D
Acq On : 14 Mar 2019 15:03
Operator : JU/SJ
Sample : K1866-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
MR-MH-04-COMP

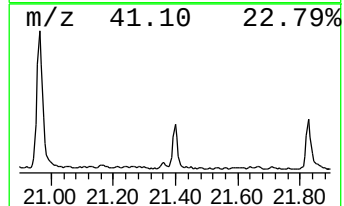
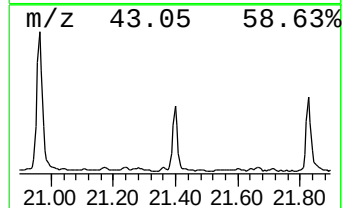
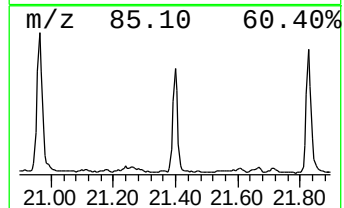
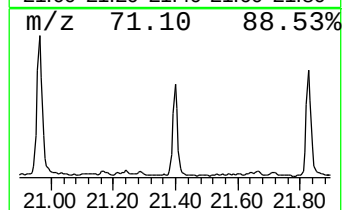
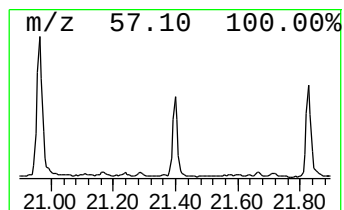
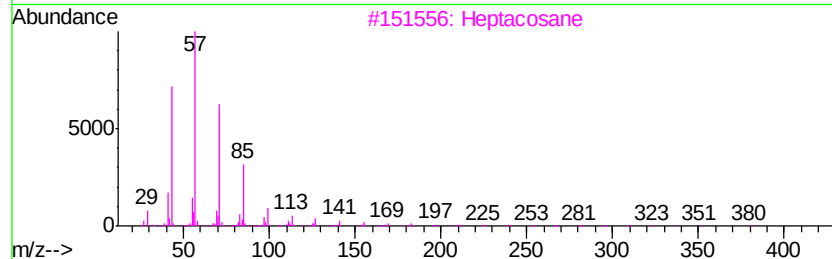
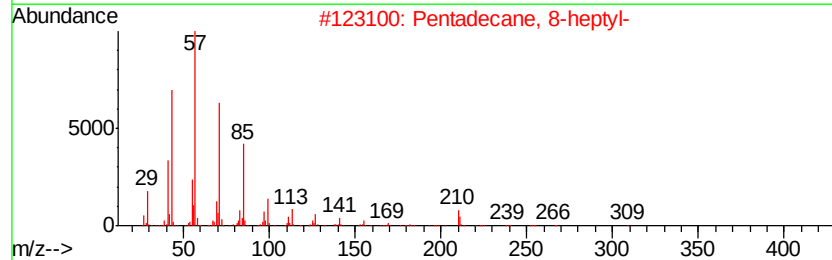
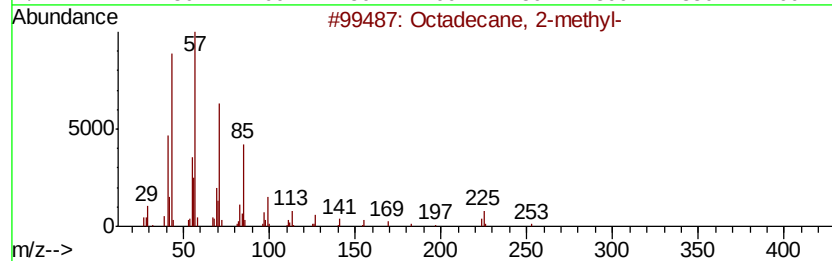
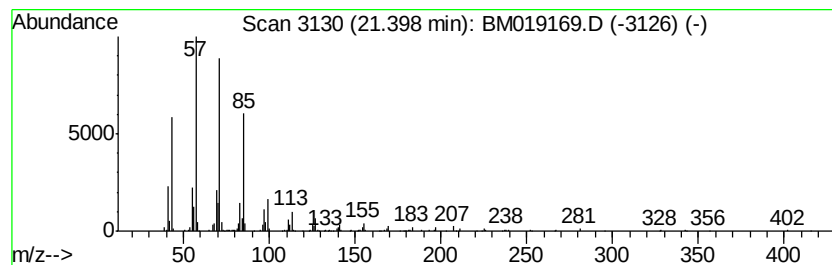
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Octadecane, 2-methyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.40	2.09 ng	403435	Chrysene-d12	21.26

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecane, 2-methyl-	268	C19H40	001560-88-9	90
2		Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	90
3		Heptacosane	380	C27H56	000593-49-7	87
4		Hexadecane	226	C16H34	000544-76-3	86
5		Nonadecane	268	C19H40	000629-92-5	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM031419\
Data File : BM019169.D
Acq On : 14 Mar 2019 15:03
Operator : JU/SJ
Sample : K1866-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

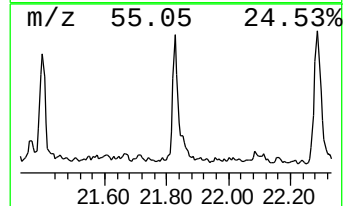
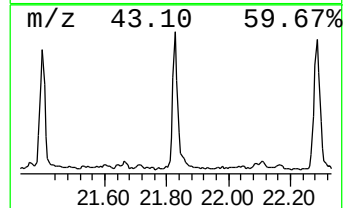
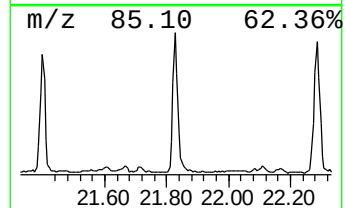
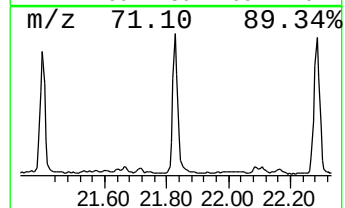
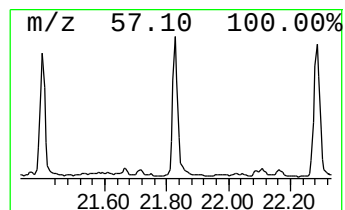
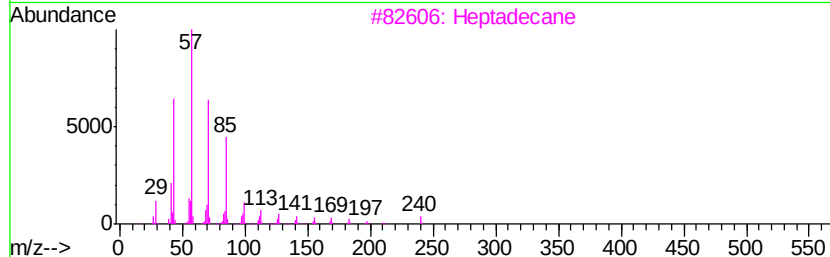
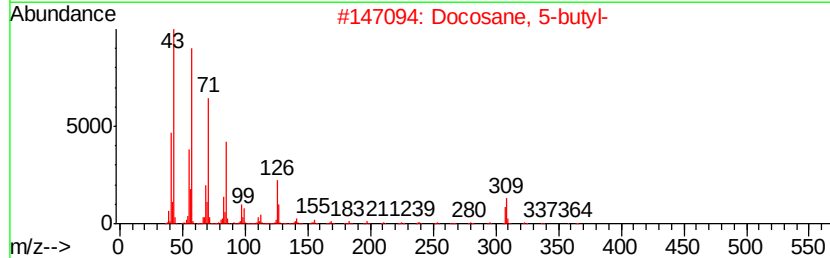
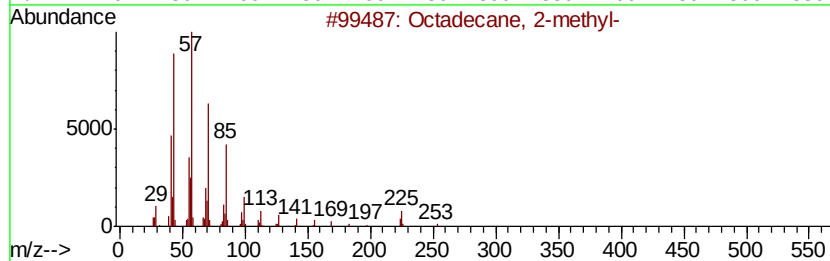
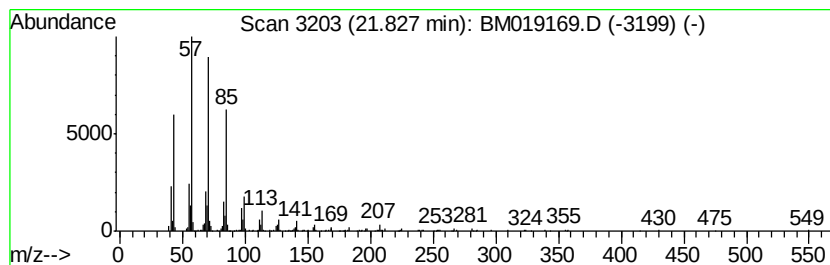
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM031119.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Docosane, 5-butyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.83	2.78 ng	536048	Chrysene-d12	21.26
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Octadecane, 2-methyl-	268 C19H40	001560-88-9 90
2		Docosane, 5-butyl-	366 C26H54	055282-16-1 90
3		Heptadecane	240 C17H36	000629-78-7 86
4		Hexadecane	226 C16H34	000544-76-3 83
5		Octadecane	254 C18H38	000593-45-3 83



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM031419\
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Acq On : 14 Mar 2019 15:03
Operator : JU/SJ
Sample : K1866-01
Misc :
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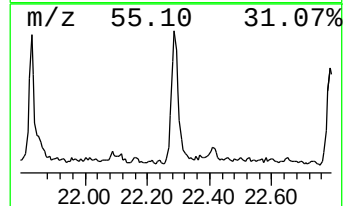
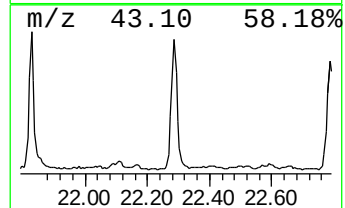
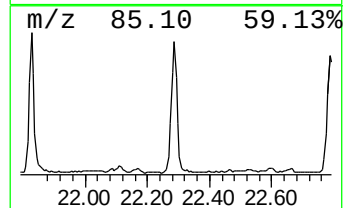
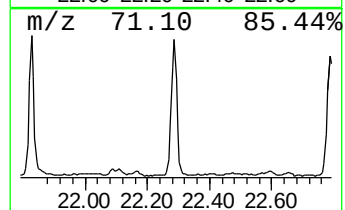
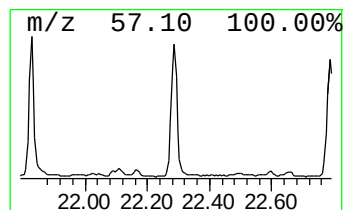
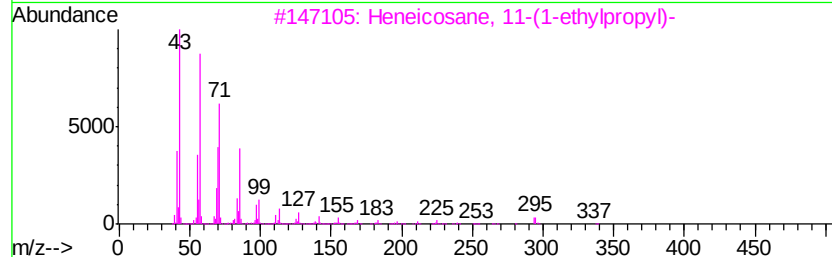
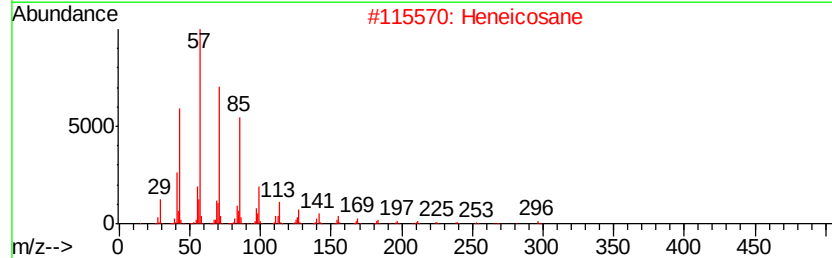
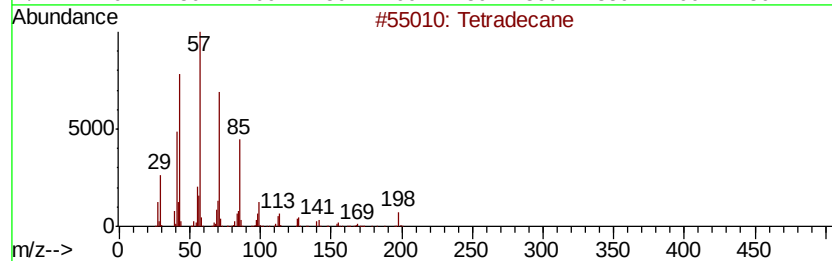
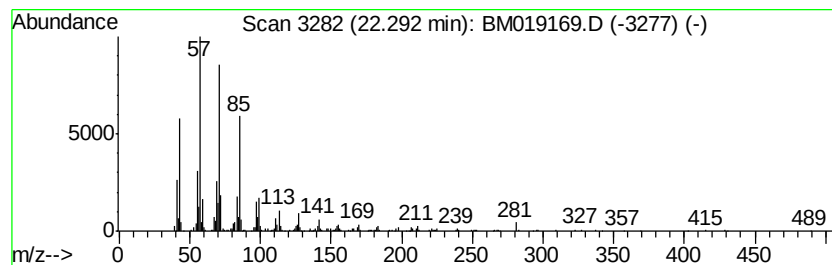
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Tetradecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD		R.T.
22.29	3.07 ng	591640	Chrysene-d12		21.26
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetradecane	198	C14H30	000629-59-4	90
2	Heneicosane	296	C21H44	000629-94-7	90
3	Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	87
4	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	87
5	Nonacosane	408	C29H60	000630-03-5	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM031419\
Data File : BM019169.D
Acq On : 14 Mar 2019 15:03
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Misc :
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Instrument :
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ClientSampleId :
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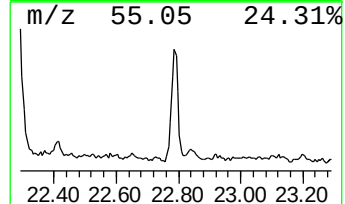
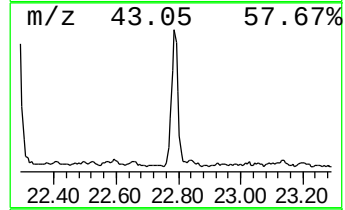
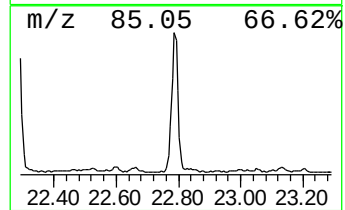
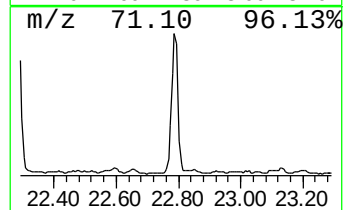
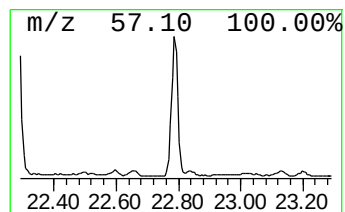
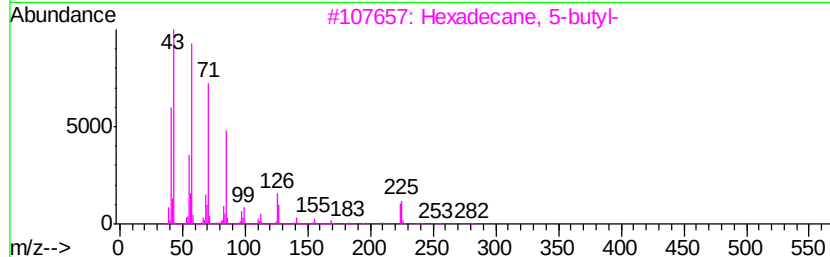
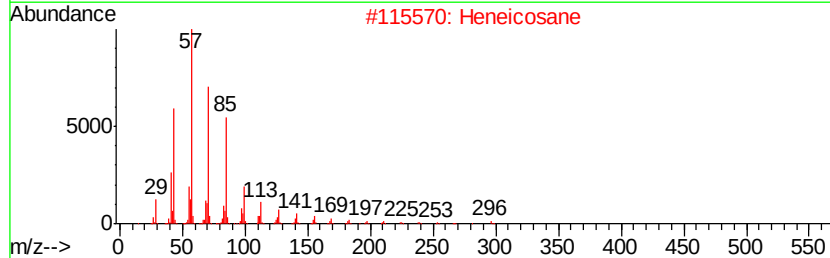
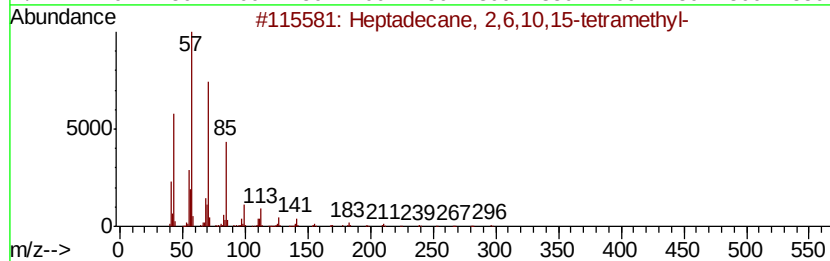
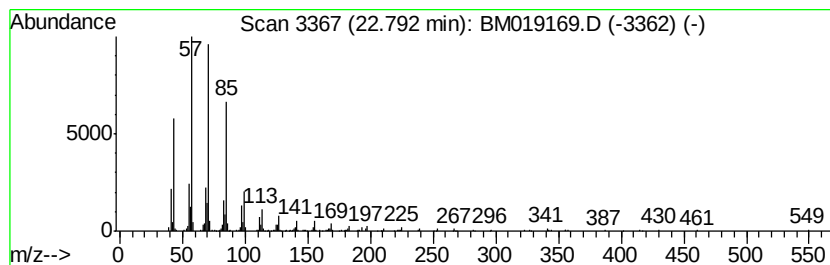
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 Heptadecane, 2,6,10,15-tetr... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.79	2.88 ng	465200	Perylene-d12	23.50

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	93
2			Heneicosane	296	C21H44	000629-94-7	91
3			Hexadecane, 5-butyl-	282	C20H42	006912-07-8	90
4			Docosane	310	C22H46	000629-97-0	87
5			Octadecane	254	C18H38	000593-45-3	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM031419\
 Data File : BM019169.D
 Acq On : 14 Mar 2019 15:03
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 Misc :
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Instrument :
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 ClientSampleId :
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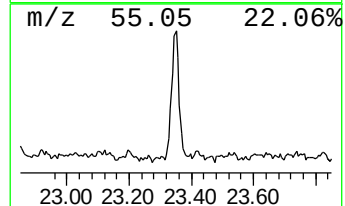
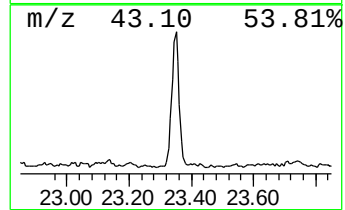
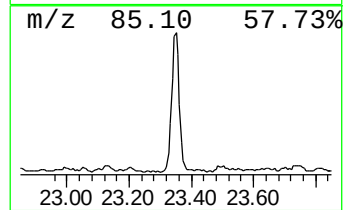
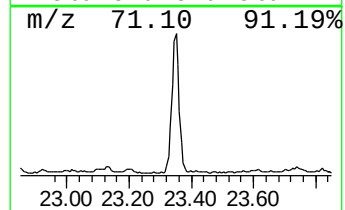
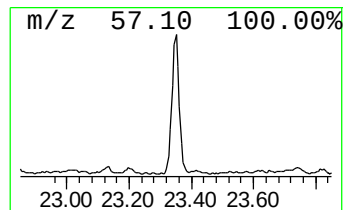
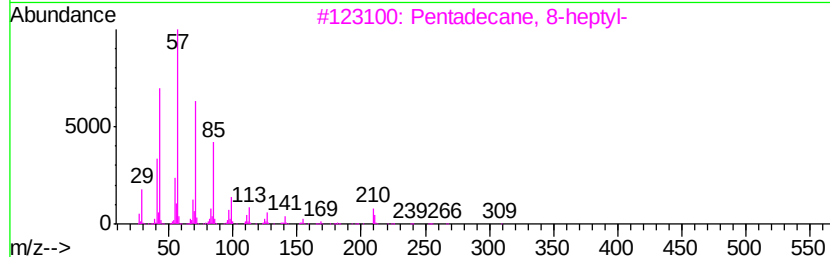
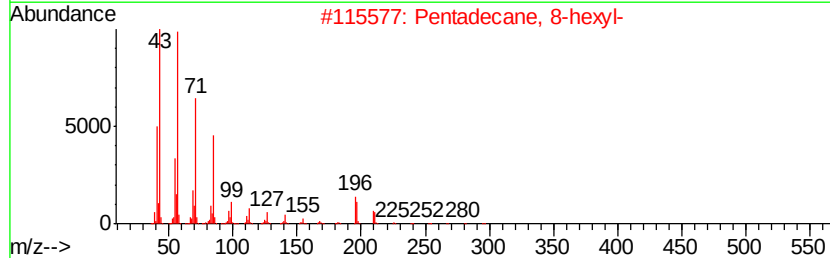
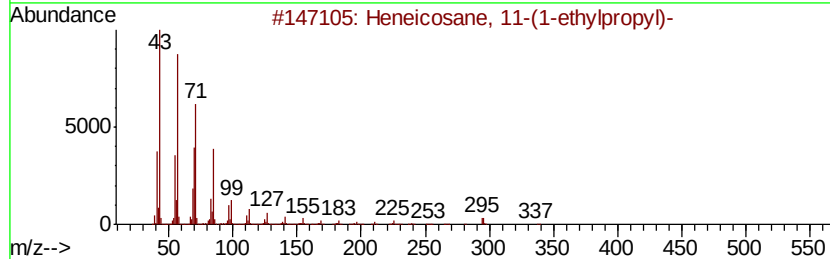
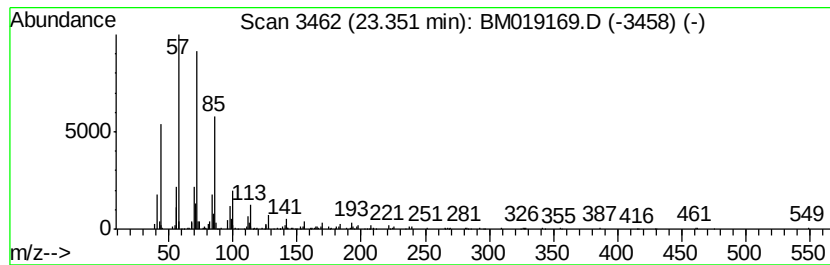
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
 TIC Integration Parameters: LSCINT.P

 Peak Number 12 Heneicosane, 11-(1-ethylpro... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.35	2.75 ng	443454	Perylene-d12	23.50

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	91
2		Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	90
3		Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	90
4		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	90
5		Heptacosane	380	C27H56	000593-49-7	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM031419\
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Acq On : 14 Mar 2019 15:03
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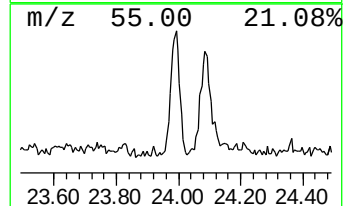
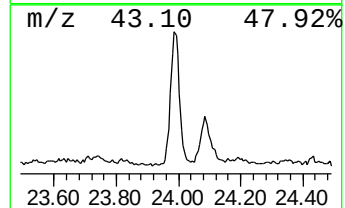
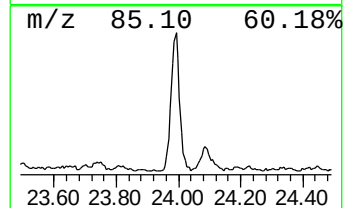
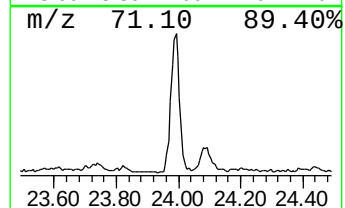
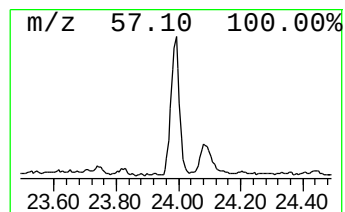
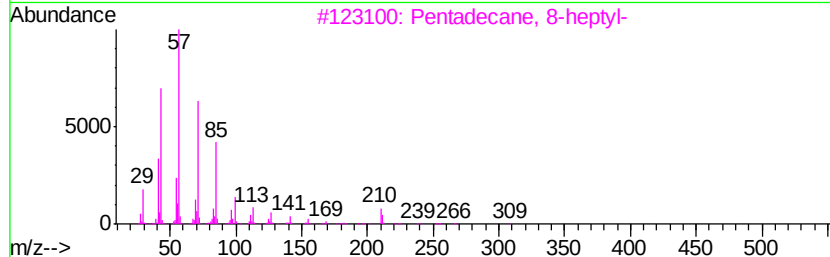
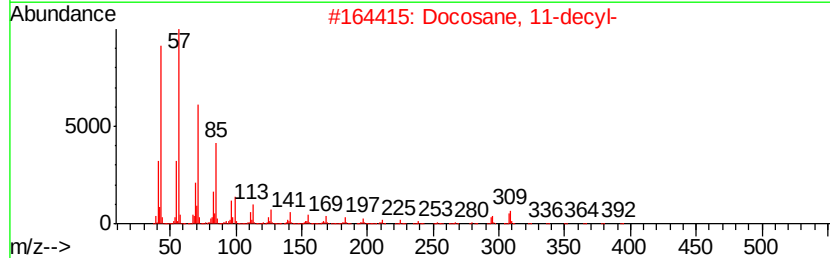
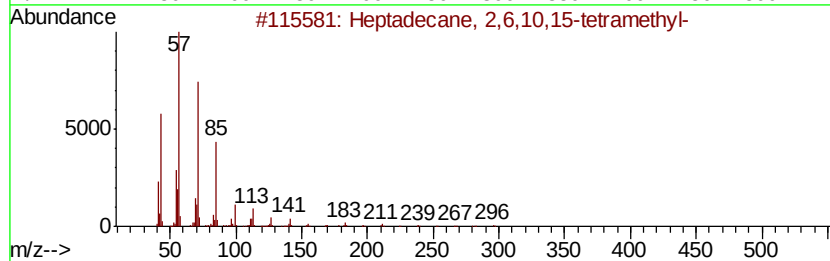
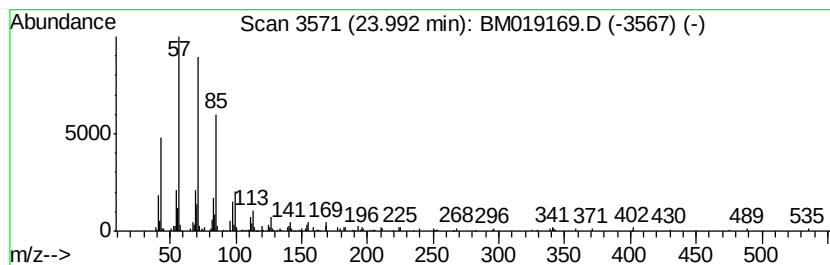
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 Heptadecane, 2,6,10,15-tetr... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.99	2.19 ng	352961	Perylene-d12	23.50

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	91
2		Docosane, 11-decyl-	451	C32H66	055401-55-3	91
3		Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	90
4		Tetracosane, 11-decyl-	479	C34H70	055429-84-0	90
5		Octadecane, 1-iodo-	380	C18H37I	000629-93-6	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM031419\
 Data File : BM019169.D
 Acq On : 14 Mar 2019 15:03
 Operator : JU/SJ
 Sample : K1866-01
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\8270-BM031119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown7.19	7.19	125.5	ng	7494460	1	7.66	1194830	20.0
2-Propenal, 3-(2,...	13.19	2.3	ng	296958	3	14.31	2645100	20.0
n-Hexadecanoic acid	17.96	8.1	ng	1375120	4	17.06	3414570	20.0
3-Hexadecene, (Z)-	18.82	7.7	ng	1314140	4	17.06	3414570	20.0
Octadecanoic acid	19.24	2.3	ng	449137	5	21.26	3855460	20.0
3-Octadecene, (E)-	20.96	7.0	ng	1354380	5	21.26	3855460	20.0
Octadecane, 2-met...	21.40	2.1	ng	403435	5	21.26	3855460	20.0
Docosane, 5-butyl-	21.83	2.8	ng	536048	5	21.26	3855460	20.0
Tetradecane	22.29	3.1	ng	591640	5	21.26	3855460	20.0
Heptadecane, 2,6,...	22.79	2.9	ng	465200	6	23.50	3228000	20.0
Heneicosane, 11-(...	23.35	2.8	ng	443454	6	23.50	3228000	20.0
Heptadecane, 2,6,...	23.99	2.2	ng	352961	6	23.50	3228000	20.0