

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF100318\
 Data File : BF109550.D
 Acq On : 3 Oct 2018 16:59
 Operator : JU/SJ
 Sample : J5194-03 2X
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HR-06-100218-B

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-BF092218.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	4.887	472	476	484	rBV	29132	38609	1.79%	0.253%
2	5.316	545	549	573	rBV	600074	810608	37.56%	5.312%
3	6.340	720	723	742	rBV	664520	887514	41.13%	5.816%
4	6.469	742	745	757	rVV	808797	916436	42.47%	6.006%
5	6.551	757	759	764	rVB	24568	23700	1.10%	0.155%
6	6.704	782	785	794	rBV	474100	494822	22.93%	3.243%
7	6.857	808	811	820	rBV	747652	670452	31.07%	4.394%
8	7.263	876	880	886	rBV	453531	490331	22.72%	3.213%
9	7.310	886	888	893	rVB	37938	39750	1.84%	0.260%
10	7.981	999	1002	1010	rBV	580309	621162	28.79%	4.071%
11	8.281	1049	1053	1056	rBV4	15302	21727	1.01%	0.142%
12	8.416	1072	1076	1081	rBV2	29249	33381	1.55%	0.219%
13	8.587	1102	1105	1108	rBV	61892	48126	2.23%	0.315%
14	9.016	1176	1178	1181	rVB2	33165	27499	1.27%	0.180%
15	9.057	1181	1185	1197	rBV	889171	828804	38.41%	5.431%
16	9.151	1197	1201	1204	rVV	59912	59013	2.73%	0.387%
17	9.451	1249	1252	1257	rBV	227874	219718	10.18%	1.440%
18	9.675	1288	1290	1293	rBV	64566	49825	2.31%	0.327%
19	9.734	1295	1300	1307	rBV	608351	575074	26.65%	3.769%
20	9.916	1325	1331	1333	rBV5	11929	23335	1.08%	0.153%
21	10.175	1371	1375	1382	rBV	60462	54020	2.50%	0.354%
22	10.392	1405	1412	1415	rBV	26731	35565	1.65%	0.233%
23	10.516	1429	1433	1442	rBV	459286	479038	22.20%	3.139%
24	10.645	1452	1455	1462	rBV	66964	85636	3.97%	0.561%
25	11.092	1527	1531	1533	rBV	57039	49121	2.28%	0.322%
26	11.122	1533	1536	1541	rVB	27496	25741	1.19%	0.169%
27	11.210	1548	1551	1559	rBV	581788	559821	25.94%	3.669%
28	11.516	1600	1603	1606	rBV	51972	38399	1.78%	0.252%
29	11.610	1616	1619	1625	rVB	463904	436999	20.25%	2.864%
30	11.745	1639	1642	1646	rBV2	85294	91106	4.22%	0.597%
31	11.922	1665	1672	1679	rBV	52802	75858	3.52%	0.497%
32	12.298	1732	1736	1738	rBV	2212933	2157909	100.00%	14.142%
33	12.316	1738	1739	1748	rVB	1673910	1447751	67.09%	9.488%
34	12.404	1751	1754	1758	rBV2	367198	371392	17.21%	2.434%

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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_F\METHODS\8270-BF092218.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	12.451	1760	1762	1766	rVB4	28215	35456	1.64%	0.232%
36	12.527	1773	1775	1779	rBV4	25413	27431	1.27%	0.180%
37	12.639	1790	1794	1798	rBV	87224	107305	4.97%	0.703%
38	12.675	1798	1800	1803	rVB	30114	28706	1.33%	0.188%
39	12.792	1816	1820	1828	rBV	794223	775732	35.95%	5.084%
40	12.957	1845	1848	1850	rBV2	59880	63743	2.95%	0.418%
41	13.033	1858	1861	1866	rBV2	59560	91010	4.22%	0.596%
42	13.122	1874	1876	1879	rBV	48487	47932	2.22%	0.314%
43	13.233	1892	1895	1899	rBV5	34792	45939	2.13%	0.301%
44	13.698	1972	1974	1983	rVB3	58685	91794	4.25%	0.602%
45	13.839	1994	1998	2005	rBV	700404	708948	32.85%	4.646%
46	15.245	2233	2237	2240	rBV	349161	447019	20.72%	2.929%

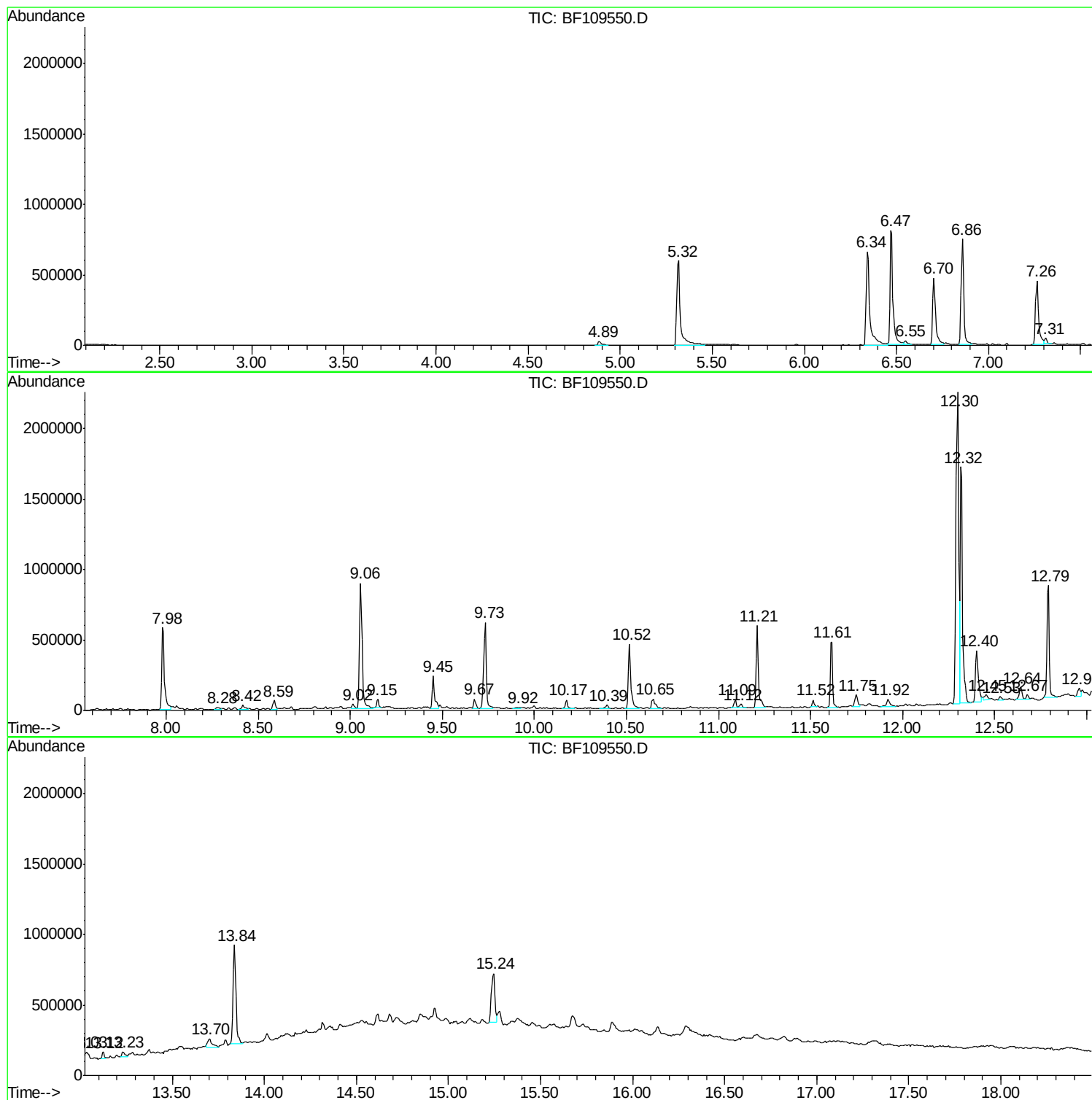
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF092218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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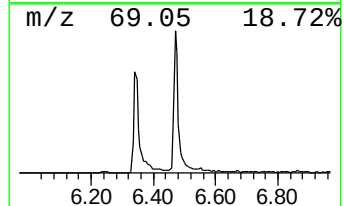
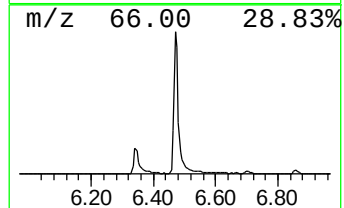
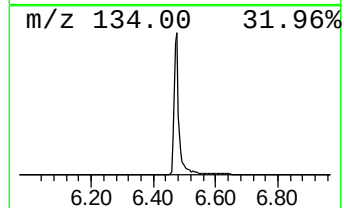
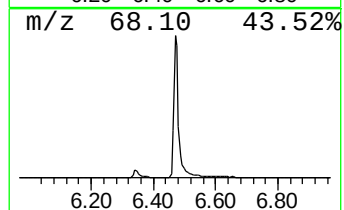
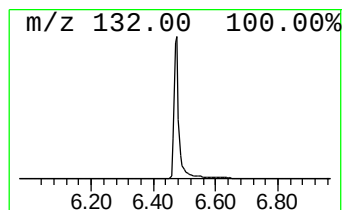
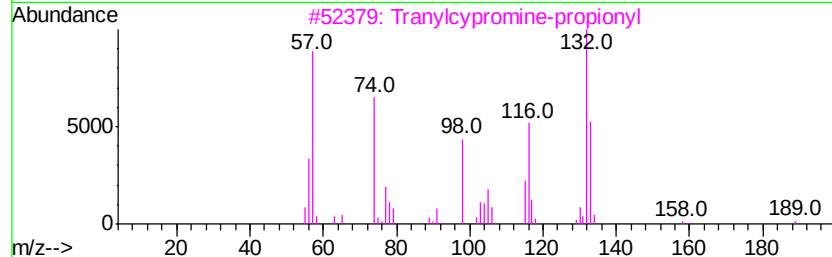
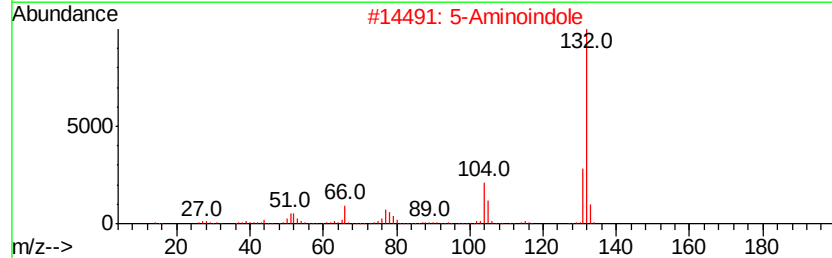
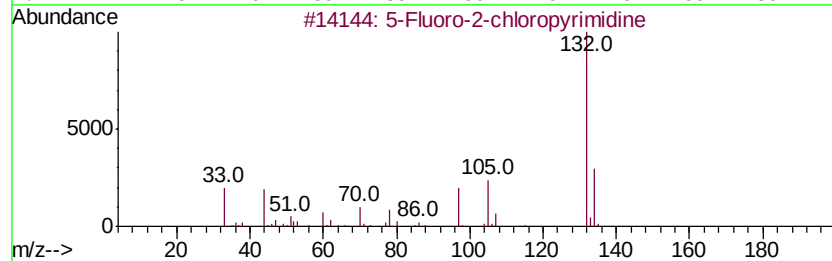
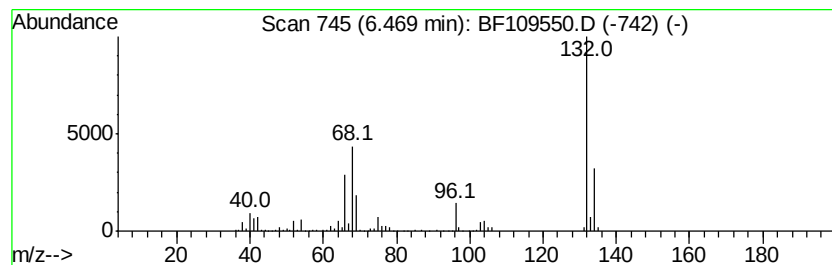
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF092218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown6.47 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.47	37.04 ng	916436	1,4-Dichlorobenzene-d4	6.70

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
2			5-Aminoindole	132	C8H8N2	005192-03-0	12
3			Tranvlcypromine-propionyl	189	C12H15NO	1000123-86-3	10
4			1,3,4-Thiadiazole-2(3H)-thione, ...	132	C3H4N2S2	029490-19-5	9
5			Imidazole, 4,5-dicyano-1-methyl-	132	C6H4N4	019485-35-9	9



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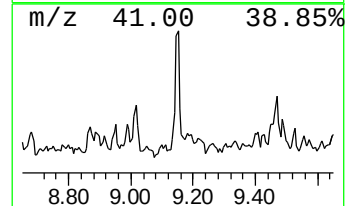
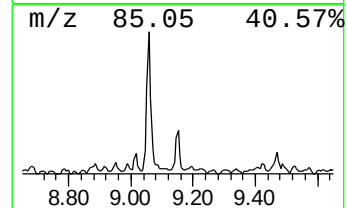
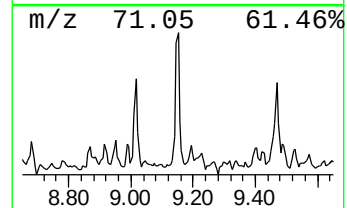
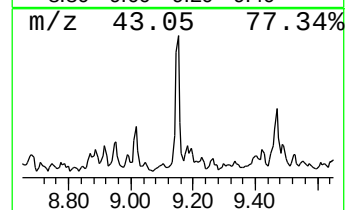
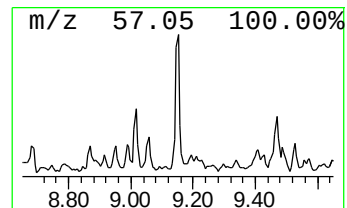
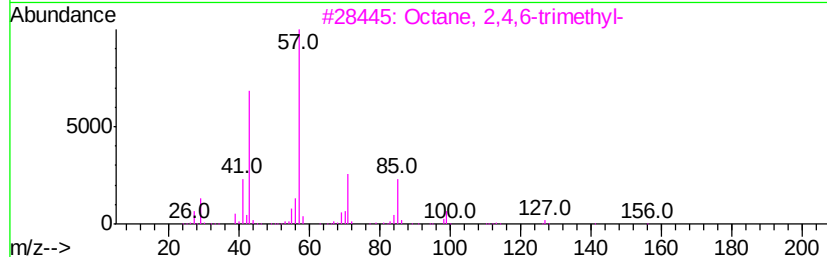
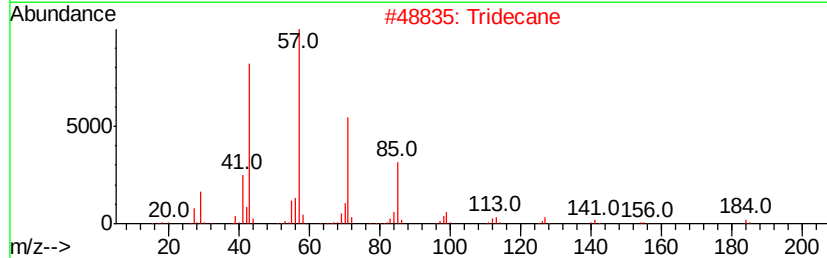
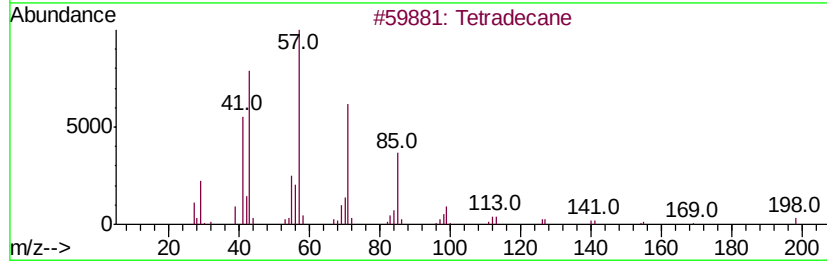
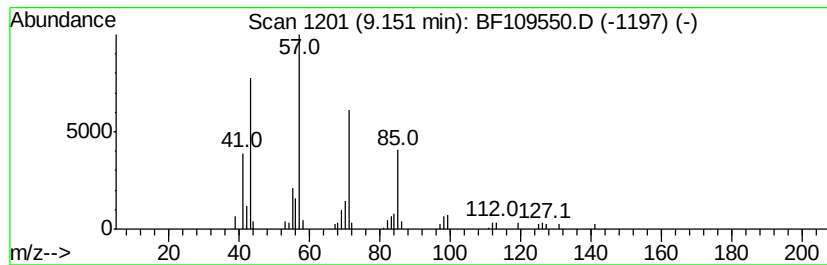
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF092218.M
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 3 Tetradecane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.15	2.05 ng	59013	Acenaphthene-d10	9.73

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetradecane	198	C14H30	000629-59-4	87
2		Tridecane	184	C13H28	000629-50-5	87
3		Octane, 2,4,6-trimethyl-	156	C11H24	062016-37-9	80
4		Hexadecane	226	C16H34	000544-76-3	80
5		Decane, 2,3,5-trimethyl-	184	C13H28	062238-11-3	78



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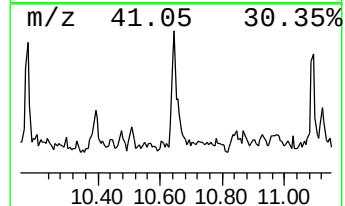
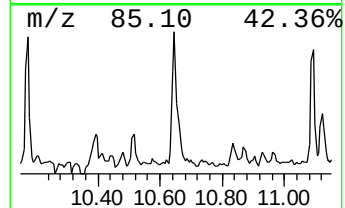
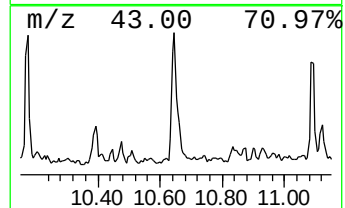
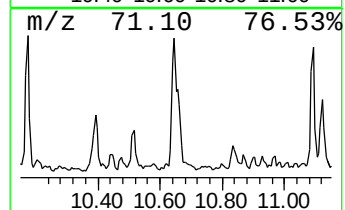
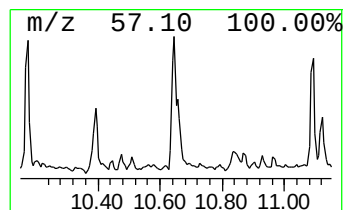
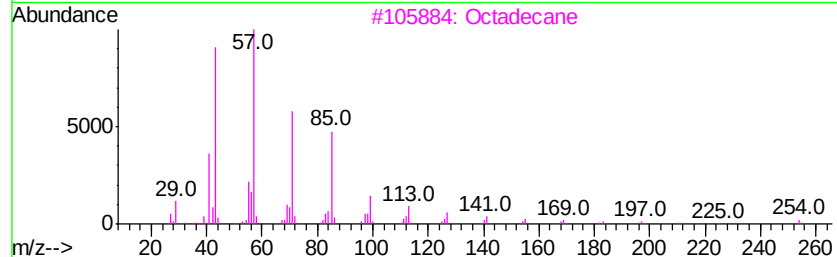
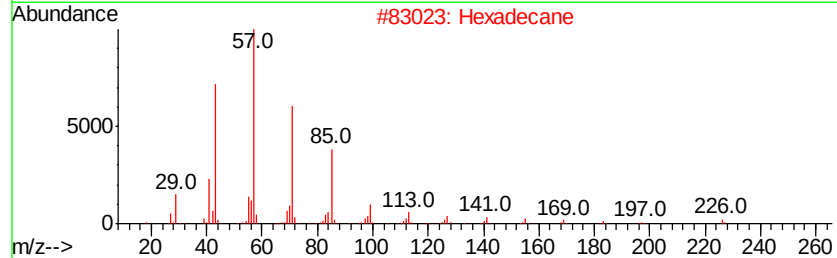
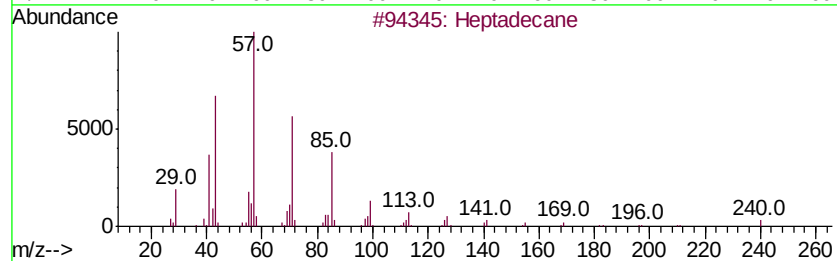
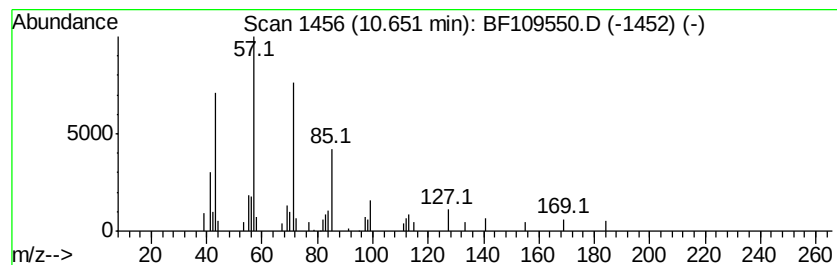
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Hexadecane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD		R.T.	
10.65	3.06 ng	85636	Phenanthrene-d10		11.21	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane		240	C17H36	000629-78-7	91
2	Hexadecane		226	C16H34	000544-76-3	91
3	Octadecane		254	C18H38	000593-45-3	90
4	Tetradecane		198	C14H30	000629-59-4	87
5	Tetracosane		338	C24H50	000646-31-1	87



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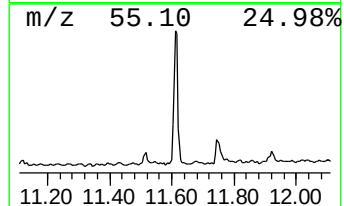
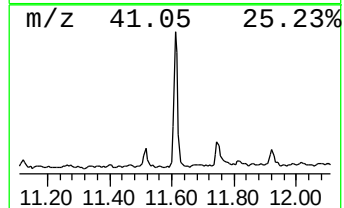
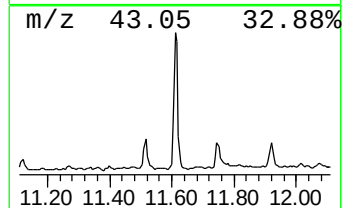
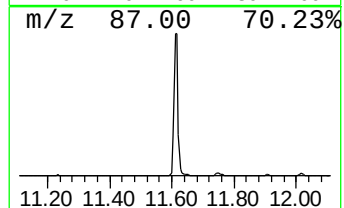
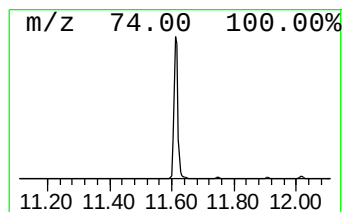
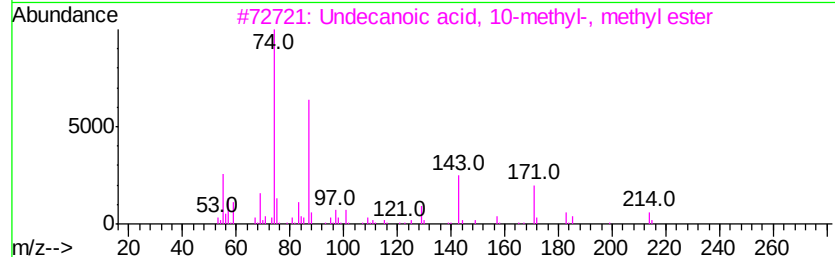
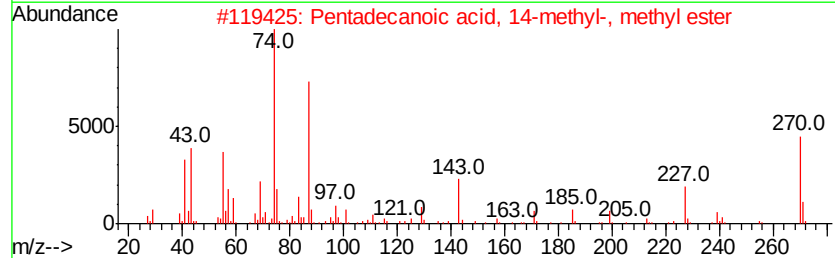
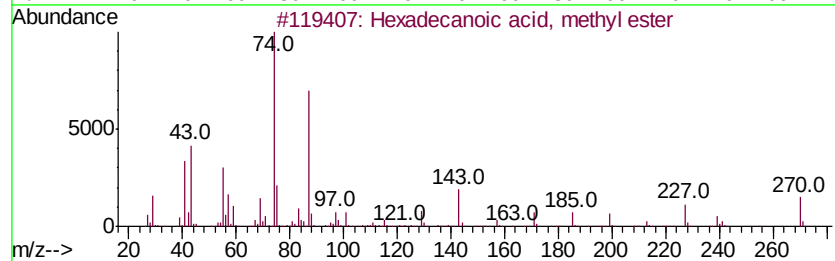
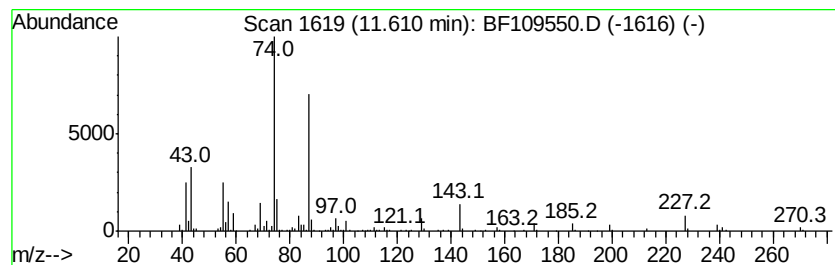
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Hexadecanoic acid, methyl e... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.61	15.61 ng	436999	Phenanthrene-d10	11.21

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	98
2		Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	93
3		Undecanoic acid, 10-methyl-, met...	214	C13H26O2	005129-56-6	90
4		Pentadecanoic acid, methyl ester	256	C16H32O2	007132-64-1	90
5		Methyl tetradecanoate	242	C15H30O2	000124-10-7	87



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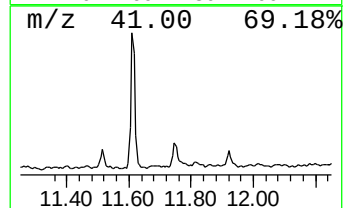
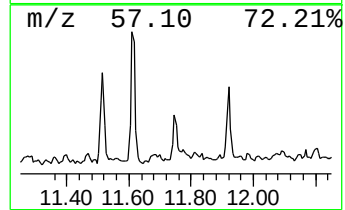
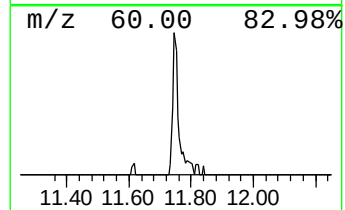
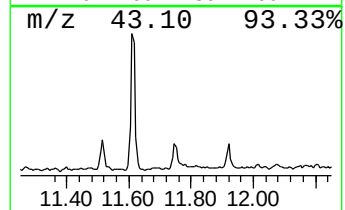
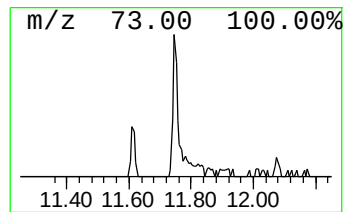
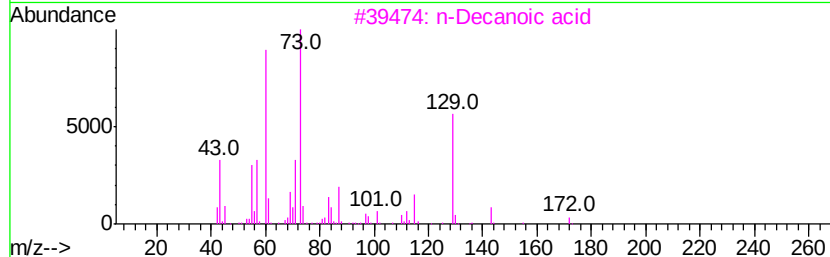
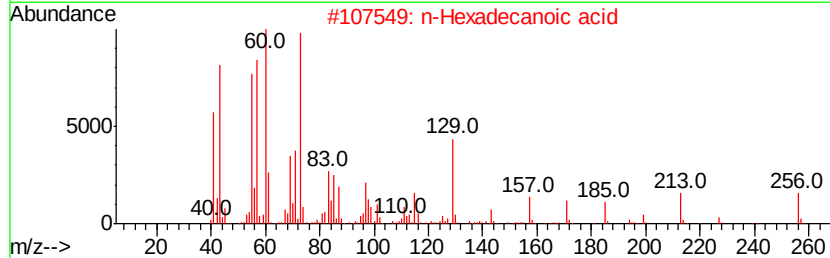
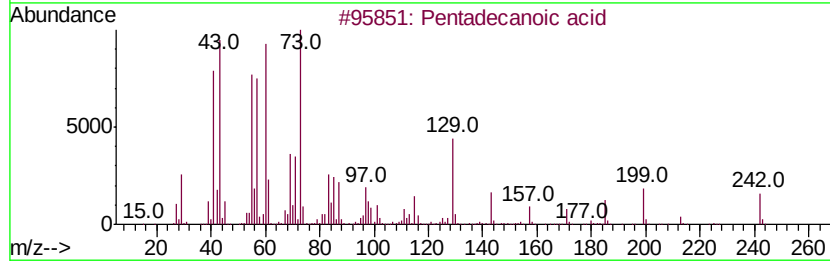
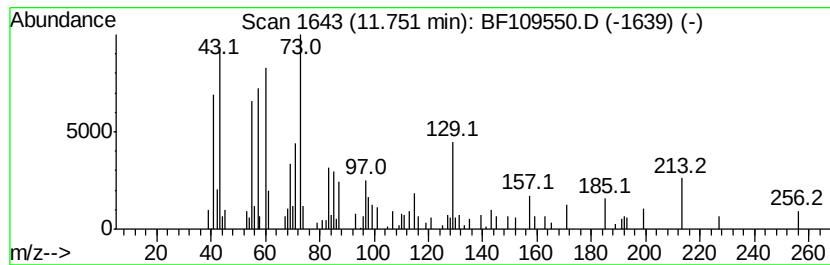
Instrument :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF092218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Pentadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD		R.T.
11.75	3.25 ng	91106	Phenanthrene-d10		11.21
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecanoic acid	242	C15H30O2	001002-84-2	68
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	64
3	n-Decanoic acid	172	C10H20O2	000334-48-5	52
4	Silane, trichlorooctadecyl-	386	C18H37Cl3Si	000112-04-9	11
5	Heneicosane	296	C21H44	000629-94-7	11



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100318\
Data File : BF109550.D
Acq On : 3 Oct 2018 16:59
Operator : JU/SJ
Sample : J5194-03 2X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
HR-06-100218-B

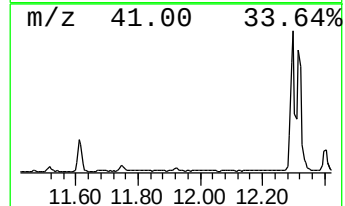
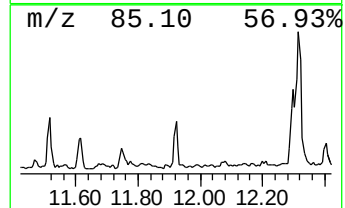
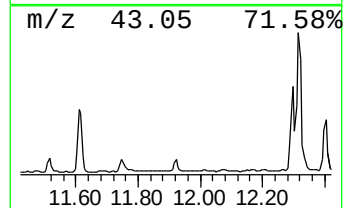
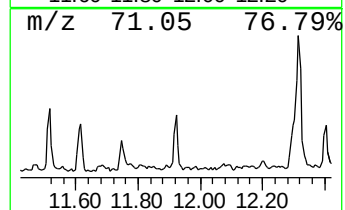
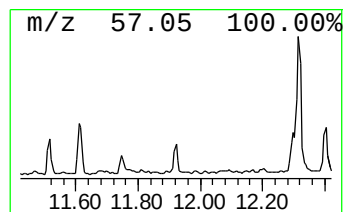
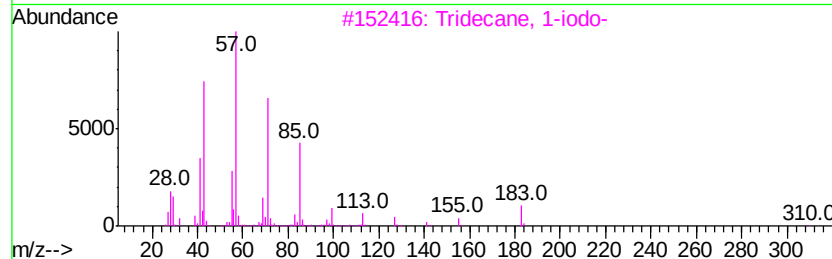
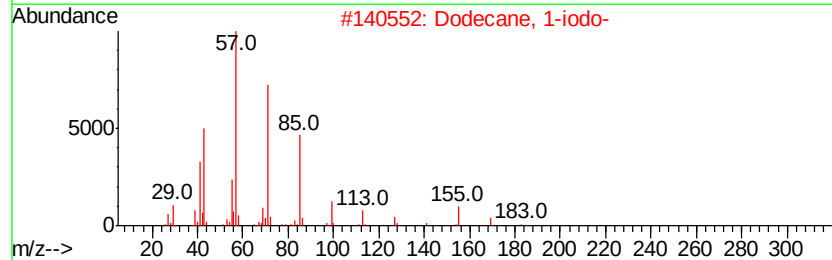
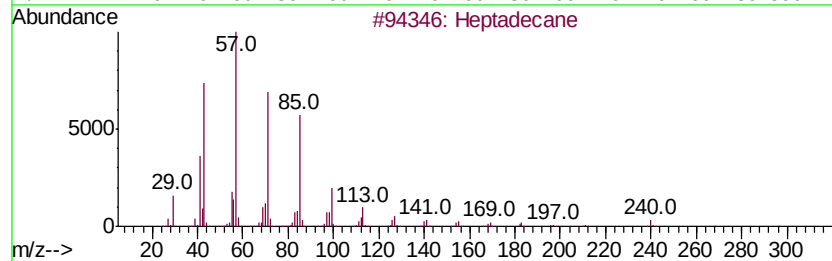
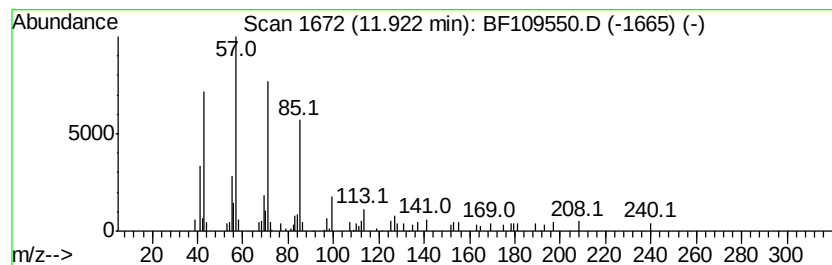
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Heptadecane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.92	2.71 ng	75858	Phenanthrene-d10	11.21

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane		240	C17H36	000629-78-7	94
2	Dodecane, 1-iodo-		296	C12H25I	004292-19-7	83
3	Tridecane, 1-iodo-		310	C13H27I	035599-77-0	83
4	Hexadecane		226	C16H34	000544-76-3	80
5	Octadecane, 1-iodo-		380	C18H37I	000629-93-6	80



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100318\
Data File : BF109550.D
Acq On : 3 Oct 2018 16:59
Operator : JU/SJ
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Misc :
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Instrument :
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ClientSampleId :
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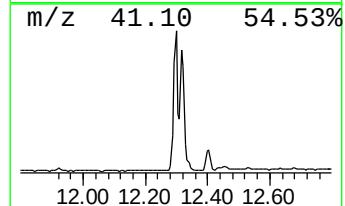
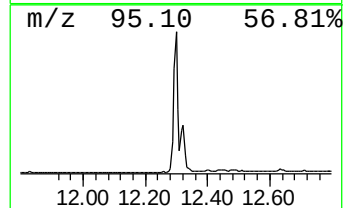
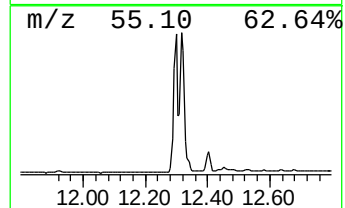
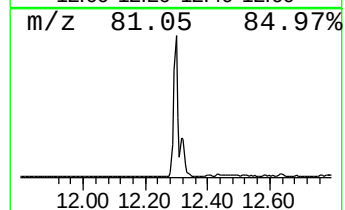
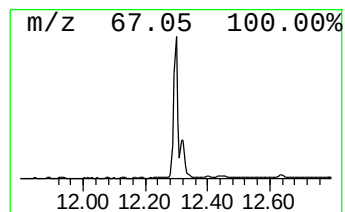
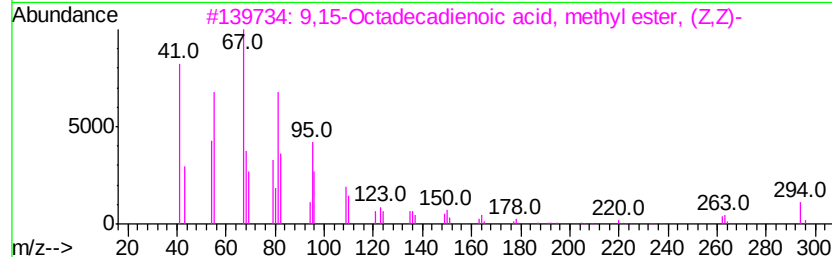
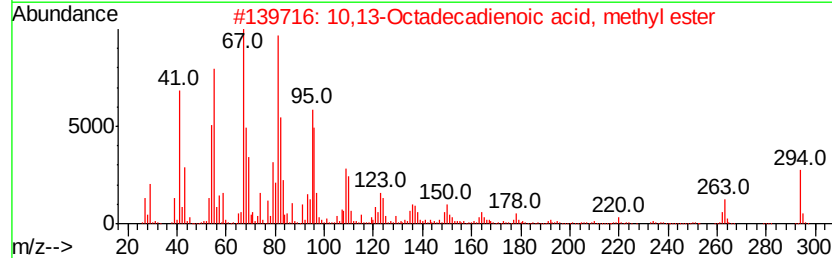
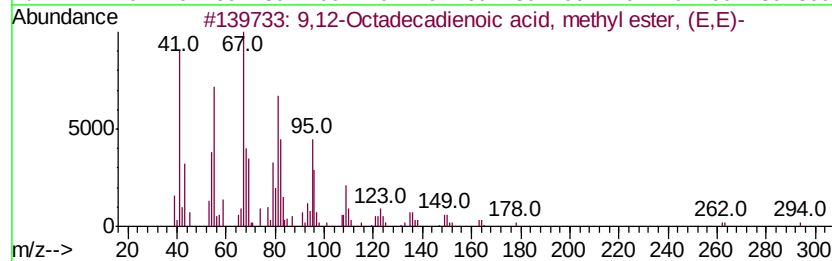
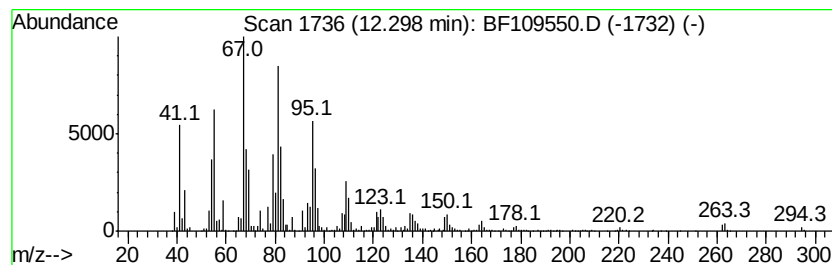
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 9,12-Octadecadienoic acid, ... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.30	77.09 ng	2157910	Phenanthrene-d10	11.21

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,12-Octadecadienoic acid, methy...	294	C19H34O2	002566-97-4	99	
2	10,13-Octadecadienoic acid, meth...	294	C19H34O2	056554-62-2	99	
3	9,15-Octadecadienoic acid, methv...	294	C19H34O2	017309-05-6	99	
4	Methyl 9-cis,11-trans-octadecadi...	294	C19H34O2	1000336-44-0	99	
5	Methyl 10-trans,12-cis-octadecad...	294	C19H34O2	1000336-44-2	98	



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100318\
Data File : BF109550.D
Acq On : 3 Oct 2018 16:59
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Sample : J5194-03 2X
Misc :
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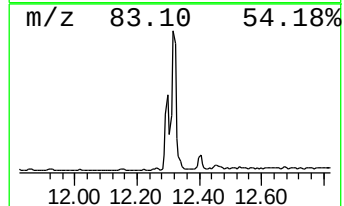
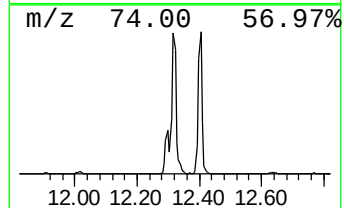
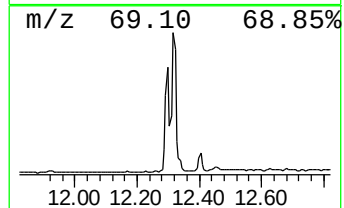
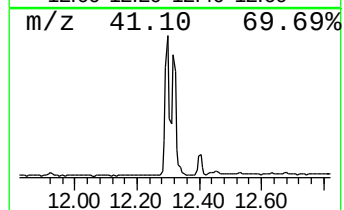
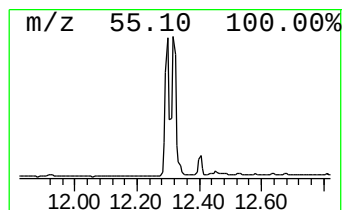
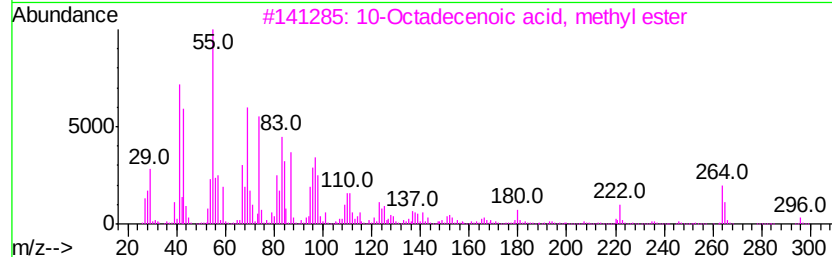
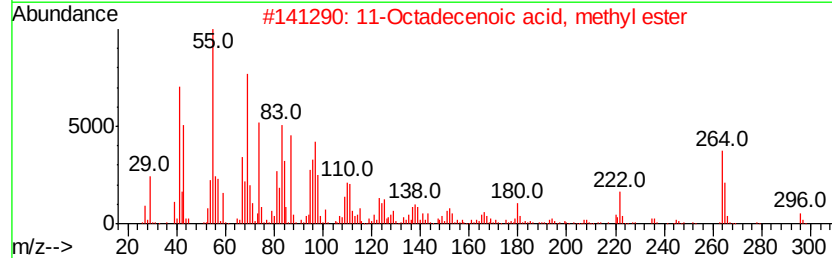
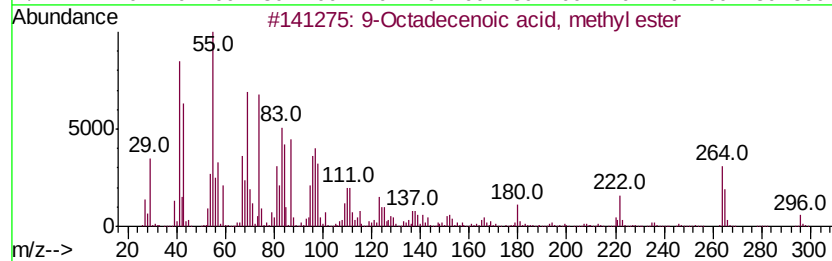
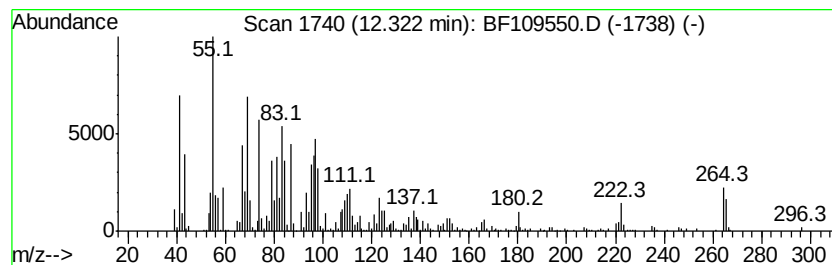
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF092218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 9-Octadecenoic acid, methyl... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.32	51.72 ng	1447750	Phenanthrene-d10	11.21	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Octadecenoic acid, methyl ester	296	C19H36O2	002462-84-2	99
2	11-Octadecenoic acid, methyl ester	296	C19H36O2	052380-33-3	99
3	10-Octadecenoic acid, methyl ester	296	C19H36O2	013481-95-3	99
4	15-Octadecenoic acid, methyl ester	296	C19H36O2	004764-72-1	99
5	9-Octadecenoic acid, methyl este...	296	C19H36O2	001937-62-8	99



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100318\
Data File : BF109550.D
Acq On : 3 Oct 2018 16:59
Operator : JU/SJ
Sample : J5194-03 2X
Misc :
ALS Vial : 12 Sample Multiplier: 1

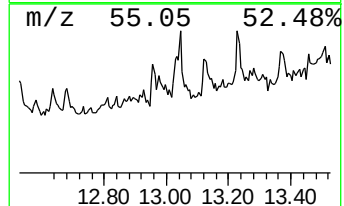
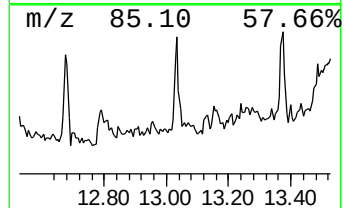
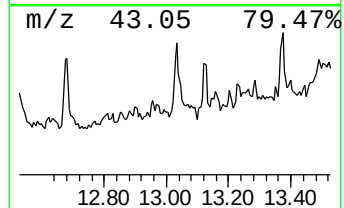
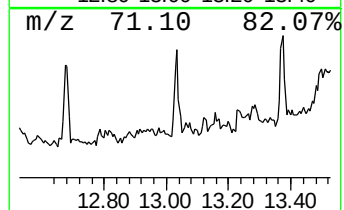
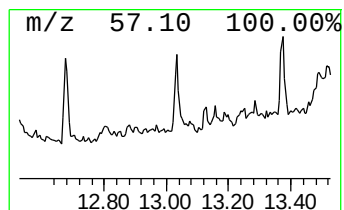
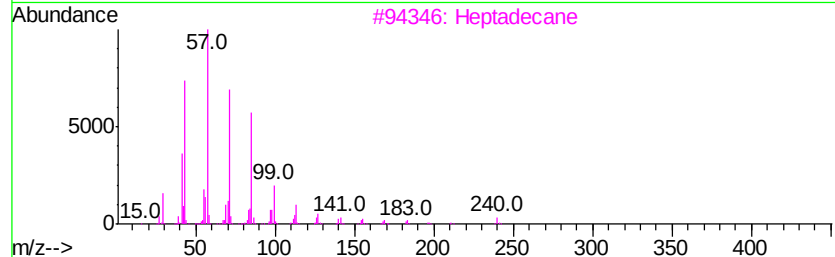
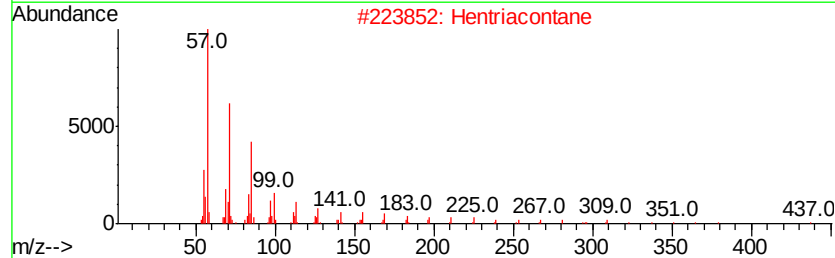
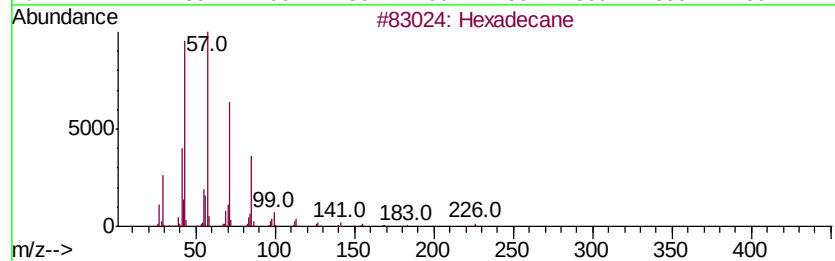
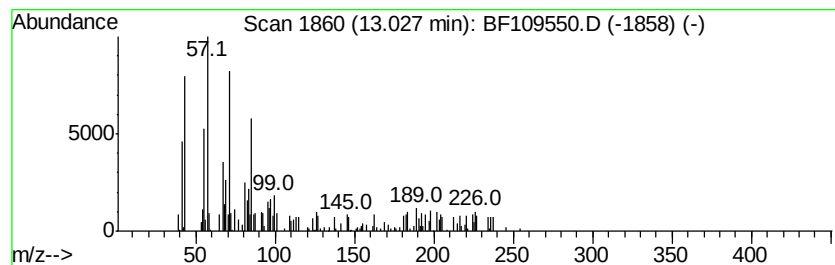
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF092218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Hentriacontane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.03	2.57 ng	91010	Chrysene-d12	13.84	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	83
2	Hentriacontane	437	C31H64	000630-04-6	59
3	Heptadecane	240	C17H36	000629-78-7	58
4	Undecane, 2,3-dimethyl-	184	C13H28	017312-77-5	52
5	Dodecane	170	C12H26	000112-40-3	52



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100318\
Data File : BF109550.D
Acq On : 3 Oct 2018 16:59
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Sample : J5194-03 2X
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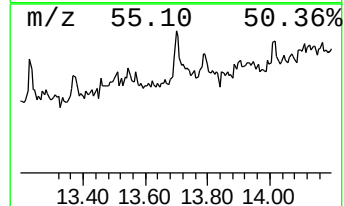
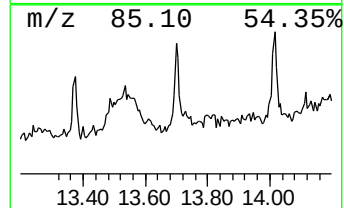
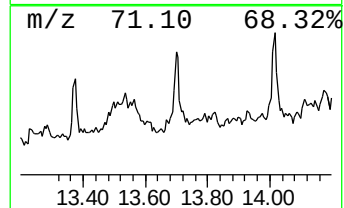
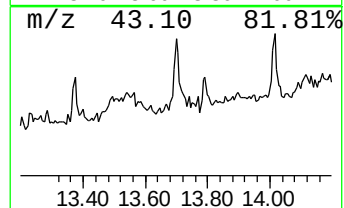
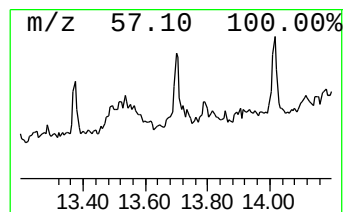
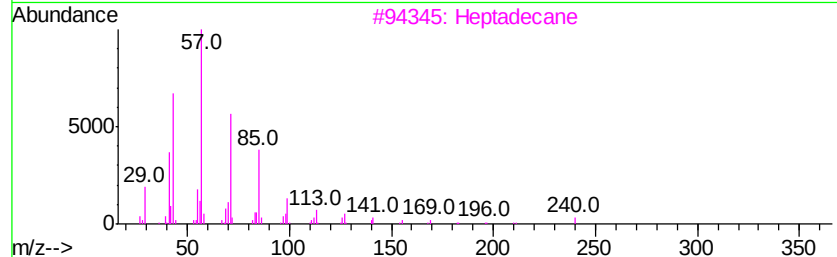
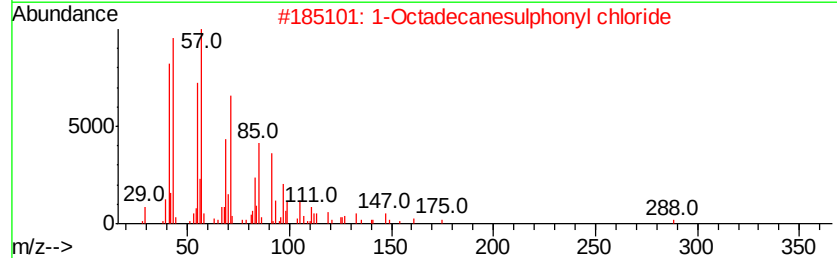
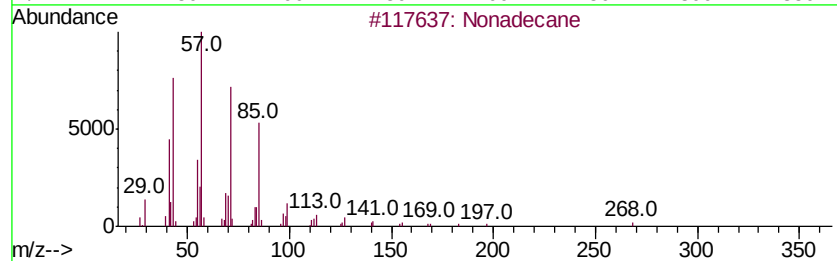
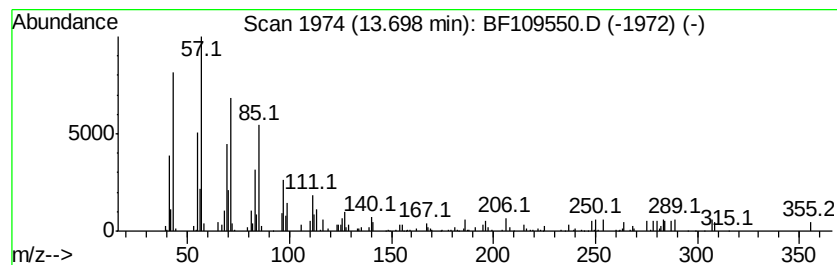
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Nonadecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD		R.T.	
13.70	2.59 ng	91794	Chrysene-d12		13.84	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Nonadecane	268	C19H40	000629-92-5	92
2		1-Octadecanesulphonyl chloride	352	C18H37ClO2S	1000342-70-4	64
3		Heptadecane	240	C17H36	000629-78-7	60
4		Octadecane	254	C18H38	000593-45-3	60
5		Eicosane	282	C20H42	000112-95-8	60



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Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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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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Hexadecane	10.65	3.1	ng	85636	4	11.21	559821	20.0
Hexadecanoic acid...	11.61	15.6	ng	436999	4	11.21	559821	20.0
Pentadecanoic acid	11.75	3.3	ng	91106	4	11.21	559821	20.0
Heptadecane	11.92	2.7	ng	75858	4	11.21	559821	20.0
9,12-Octadecadien...	12.30	77.1	ng	2157910	4	11.21	559821	20.0
9-Octadecenoic ac...	12.32	51.7	ng	1447750	4	11.21	559821	20.0
Hentriacontane	13.03	2.6	ng	91010	5	13.84	708948	20.0
Nonadecane	13.70	2.6	ng	91794	5	13.84	708948	20.0