

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072718\
 Data File : BF107762.D
 Acq On : 27 Jul 2018 22:37
 Operator : JU/SJ
 Sample : J4136-02
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CON-SB-01-02-03-0-25

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-BF072018.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	5.189	482	485	502	rBV	242379	305742	6.17%	0.800%
2	5.595	550	554	586	rBV	949987	1850406	37.36%	4.840%
3	6.595	720	724	744	rBV	1015302	1901611	38.39%	4.974%
4	6.731	744	747	781	rVV	1665106	2471989	49.91%	6.466%
5	6.960	783	786	808	rVV	496793	911876	18.41%	2.385%
6	7.113	808	812	840	rVB	1480917	1705758	34.44%	4.462%
7	7.525	878	882	904	rBV	653941	1254372	25.33%	3.281%
8	8.242	1000	1004	1022	rBV	814671	1076697	21.74%	2.817%
9	9.313	1182	1186	1201	rBV	1881636	2112023	42.64%	5.525%
10	9.413	1201	1203	1213	rVB3	30275	58409	1.18%	0.153%
11	9.707	1249	1253	1265	rBV	405633	416962	8.42%	1.091%
12	10.001	1299	1303	1306	rBV	1035439	1012372	20.44%	2.648%
13	10.030	1306	1308	1323	rVB2	106651	149632	3.02%	0.391%
14	10.795	1434	1438	1450	rBV	951695	1258482	25.41%	3.292%
15	10.989	1468	1471	1475	rVB	65229	58510	1.18%	0.153%
16	11.495	1553	1557	1559	rBV	830464	830523	16.77%	2.173%
17	11.519	1559	1561	1568	rVV	455890	570290	11.51%	1.492%
18	11.571	1568	1570	1580	rVB	103813	164152	3.31%	0.429%
19	11.789	1604	1607	1611	rVB	112724	106472	2.15%	0.279%
20	11.866	1616	1620	1627	rVV	1722413	1636427	33.04%	4.281%
21	12.001	1639	1643	1646	rBV2	175761	188123	3.80%	0.492%
22	12.113	1658	1662	1664	rBV2	77862	81403	1.64%	0.213%
23	12.554	1733	1737	1739	rBV	3798396	3989076	80.54%	10.435%
24	12.577	1739	1741	1748	rVV2	5227116	4952820	100.00%	12.956%
25	12.654	1751	1754	1760	rVV	928390	823396	16.62%	2.154%
26	12.713	1760	1764	1775	rVV	724044	801966	16.19%	2.098%
27	12.871	1788	1791	1793	rBV	84227	72780	1.47%	0.190%
28	12.942	1797	1803	1809	rBV	533230	713736	14.41%	1.867%
29	13.077	1822	1826	1836	rBV	1616685	1691039	34.14%	4.424%
30	13.165	1837	1841	1847	rVB3	34934	70056	1.41%	0.183%
31	13.271	1852	1859	1863	rVV2	64150	167177	3.38%	0.437%
32	13.307	1863	1865	1868	rVB	98488	84247	1.70%	0.220%
33	13.377	1874	1877	1881	rBV2	96389	97396	1.97%	0.255%
34	13.895	1962	1965	1967	rBV3	52717	59160	1.19%	0.155%

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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-BF072018.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	13.954	1972	1975	1980	rVV3	213541	280917	5.67%	0.735%
36	13.995	1980	1982	1987	rVB2	61834	76251	1.54%	0.199%
37	14.048	1989	1991	1997	rBV2	60155	70342	1.42%	0.184%
38	14.148	2002	2008	2011	rBV2	950413	1406288	28.39%	3.679%
39	14.171	2011	2012	2020	rVB	311319	274628	5.54%	0.718%
40	14.577	2079	2081	2084	rVB	72325	65487	1.32%	0.171%
41	14.748	2107	2110	2113	rVB	108745	112323	2.27%	0.294%
42	14.895	2132	2135	2139	rVB3	94551	104253	2.10%	0.273%
43	14.977	2145	2149	2157	rVB	518341	537605	10.85%	1.406%
44	15.218	2187	2190	2193	rBV	155861	244716	4.94%	0.640%
45	15.248	2193	2195	2202	rVB2	170824	194551	3.93%	0.509%
46	15.542	2243	2245	2252	rVB	75997	101239	2.04%	0.265%
47	15.612	2253	2257	2260	rVV	126471	181768	3.67%	0.475%
48	15.677	2264	2268	2284	rVB	522706	932300	18.82%	2.439%

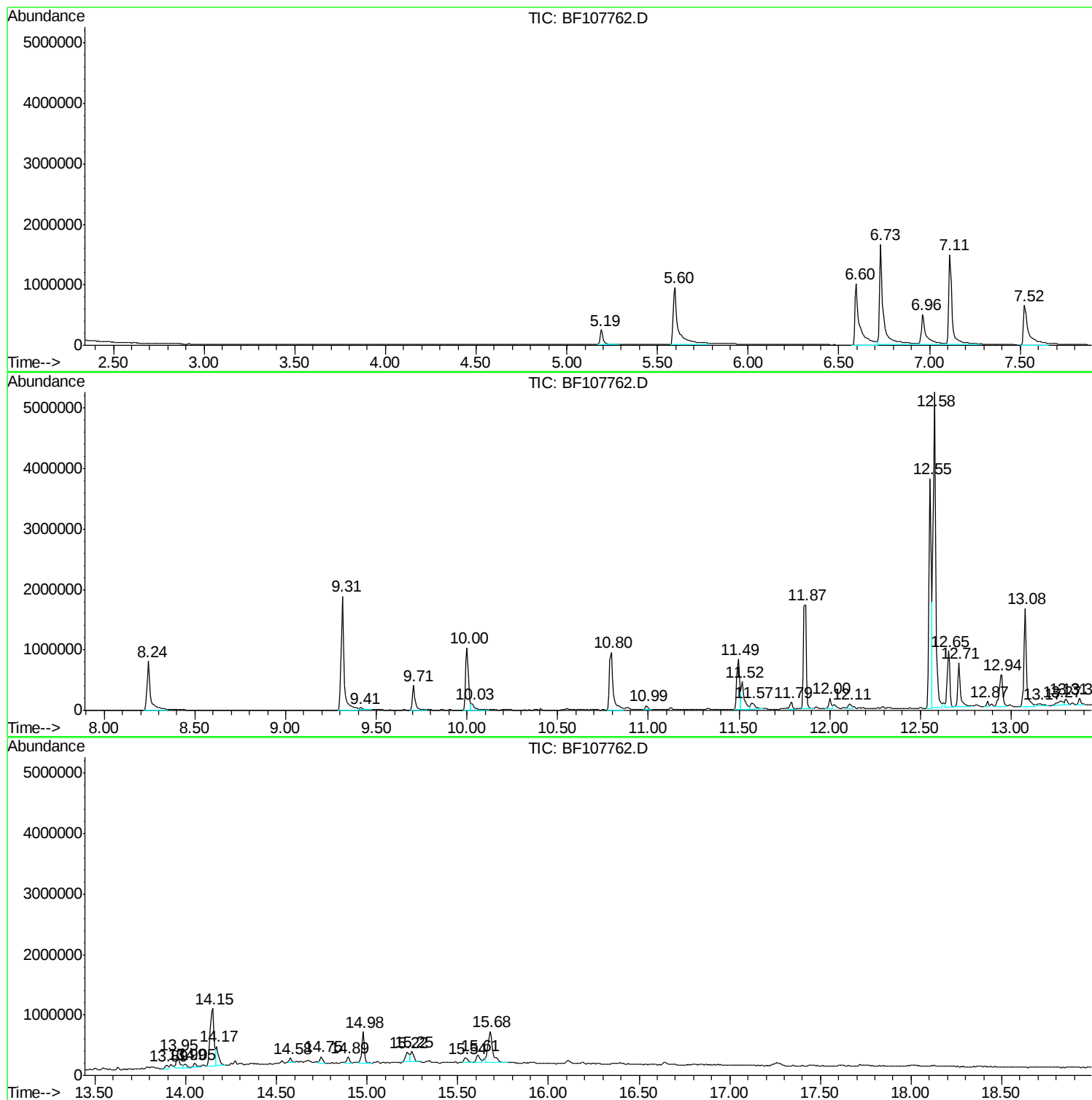
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF072018.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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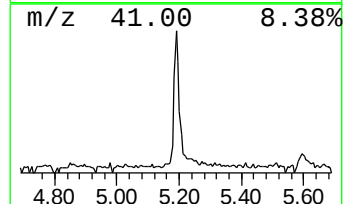
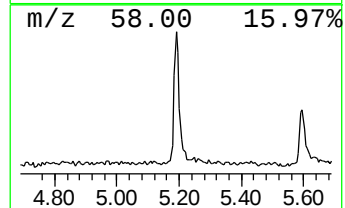
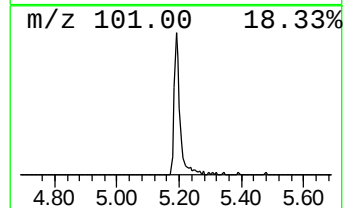
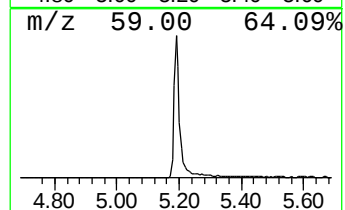
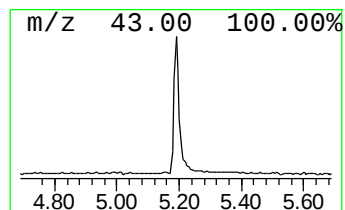
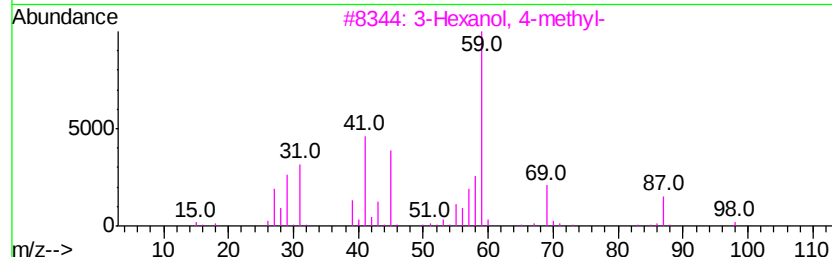
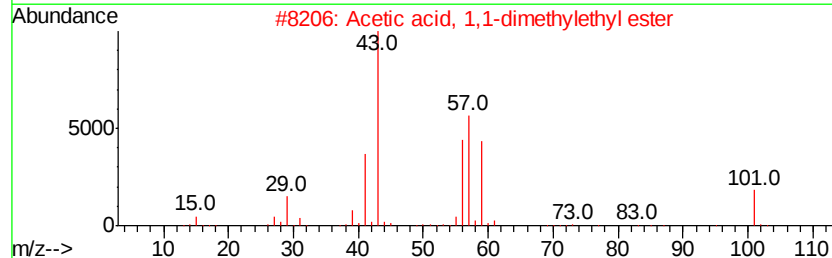
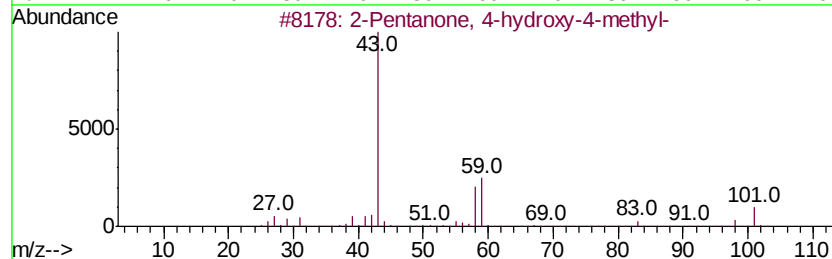
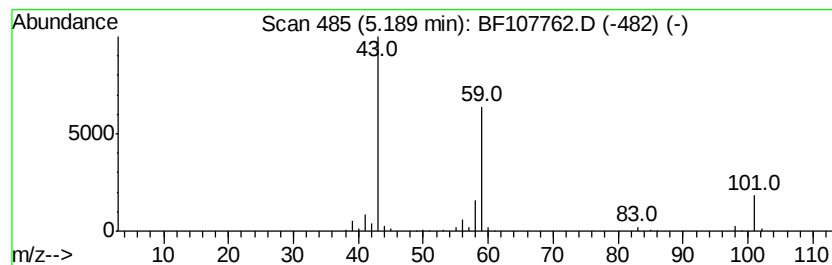
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.19	6.71 ng	305742	1,4-Dichlorobenzene-d4	6.96

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
4			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25
5			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9



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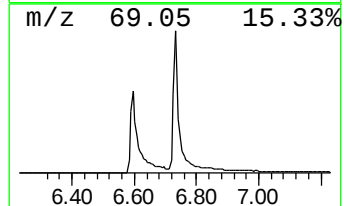
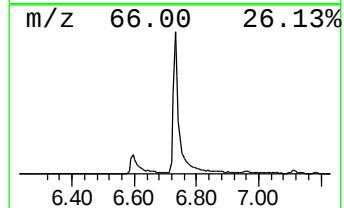
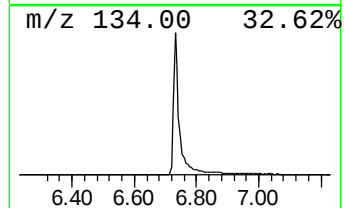
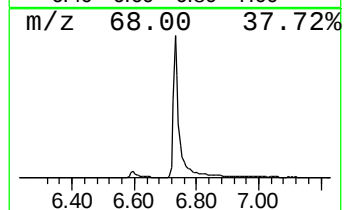
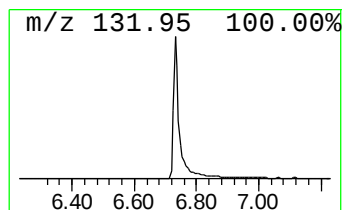
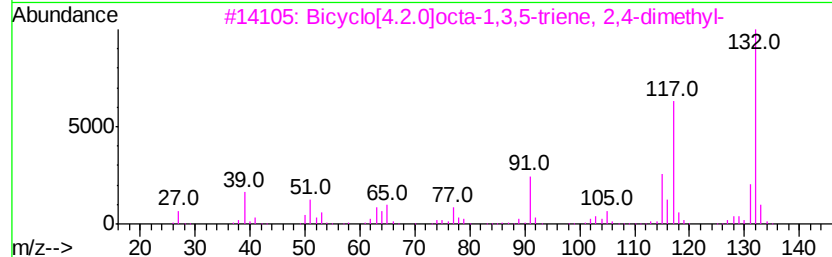
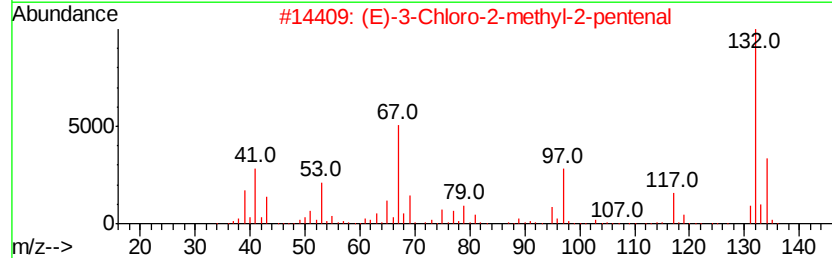
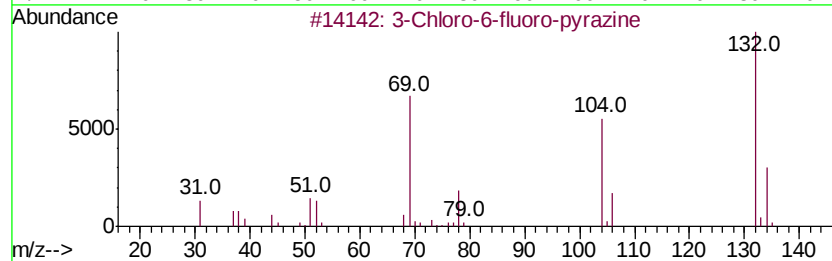
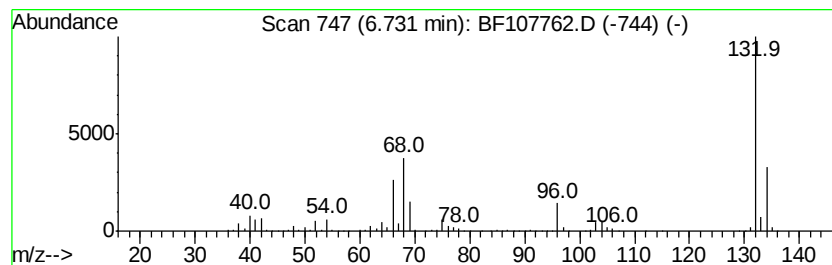
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown6.73 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.73	54.22 ng	2471990	1,4-Dichlorobenzene-d4	6.96

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9
5		4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9



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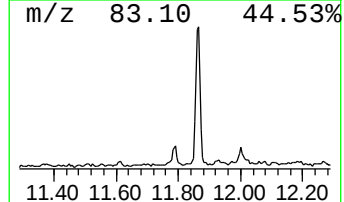
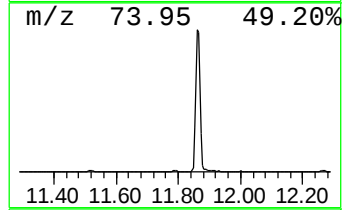
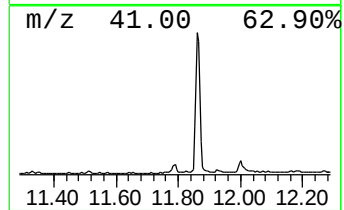
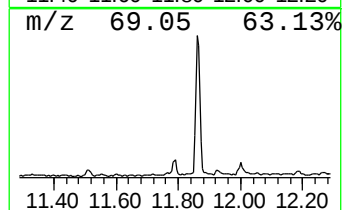
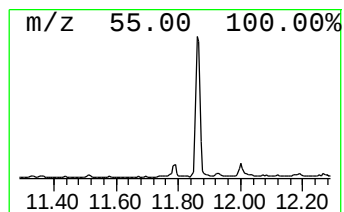
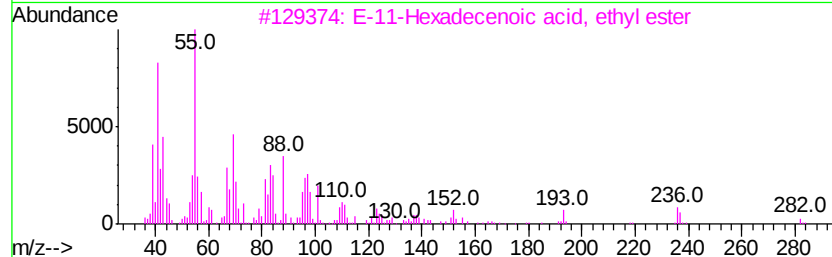
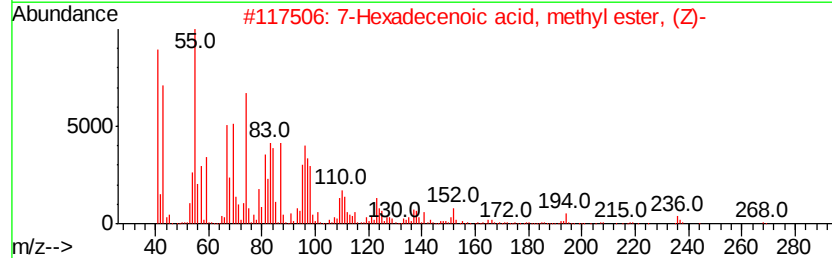
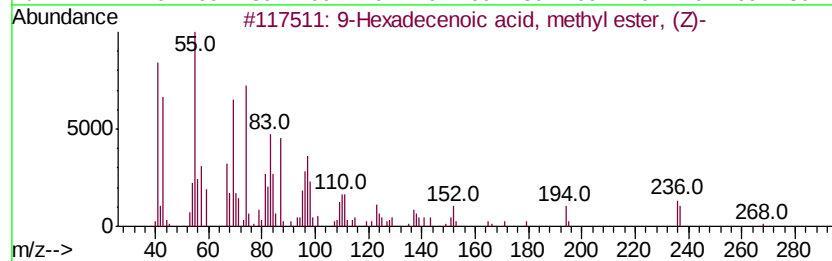
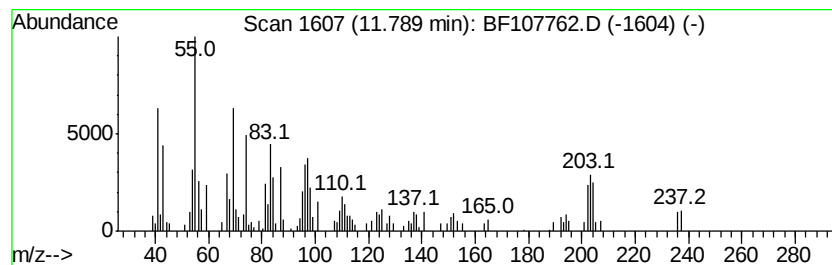
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 9-Hexadecenoic acid, methyl... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.		
11.79	2.56 ng	106472	Phenanthrene-d10	11.49		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		9-Hexadecenoic acid, methyl este...	268	C17H32O2	001120-25-8	60
2		7-Hexadecenoic acid, methyl este...	268	C17H32O2	056875-67-3	49
3		E-11-Hexadecenoic acid, ethyl ester	282	C18H34O2	1000245-71-9	43
4		Oxiraneundecanoic acid, 3-pentyl...	312	C19H36O3	038520-31-9	43
5		Cyclopentadecanone, 2-hydroxy-	240	C15H28O2	004727-18-8	41



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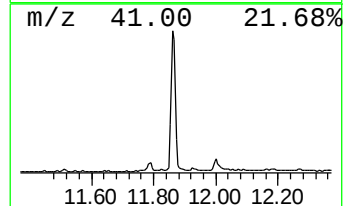
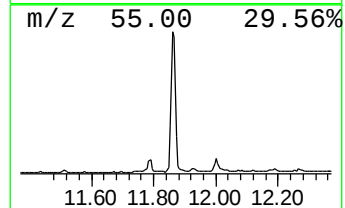
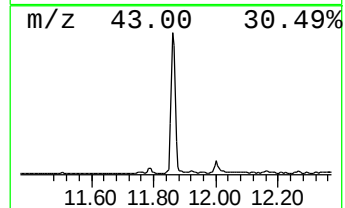
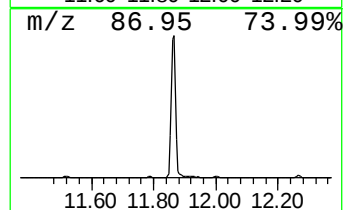
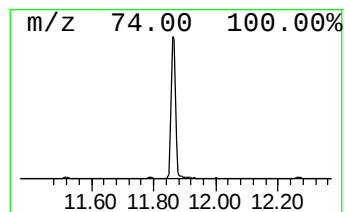
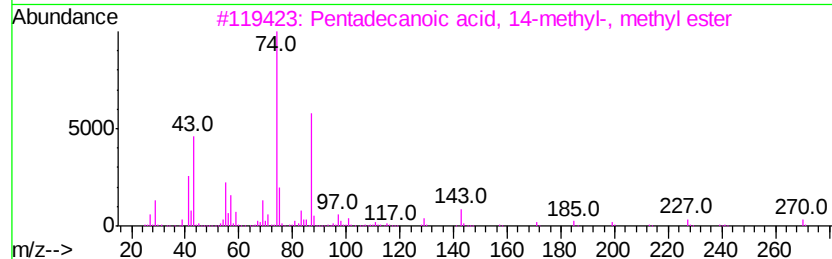
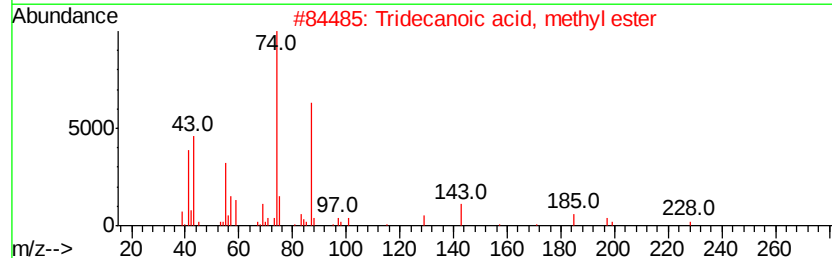
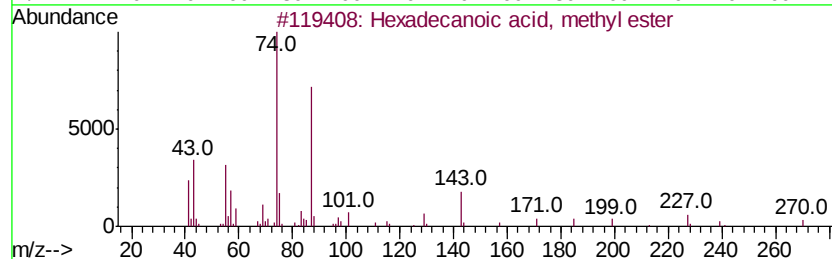
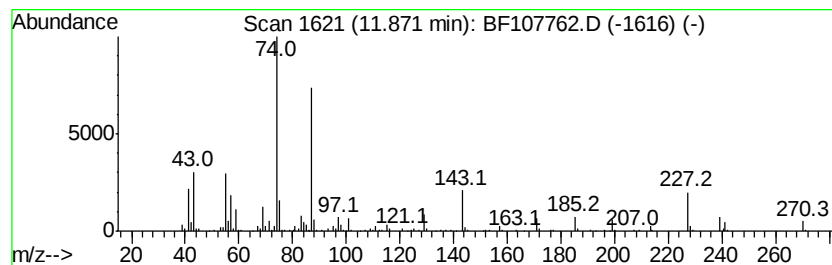
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Hexadecanoic acid, methyl e... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.87	39.41 ng	1636430	Phenanthrene-d10	11.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	98
2		Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	93
3		Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	92
4		Methyl tetradecanoate	242	C15H30O2	000124-10-7	87
5		Decanoic acid, methyl ester	186	C11H22O2	000110-42-9	87



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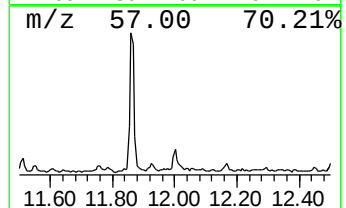
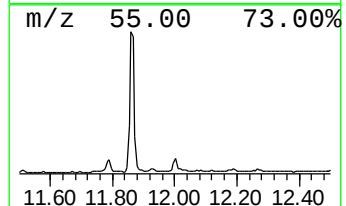
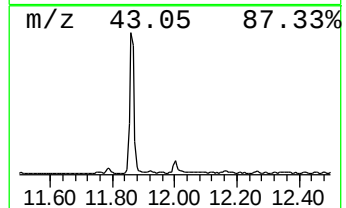
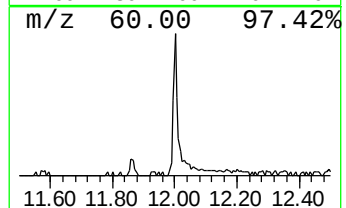
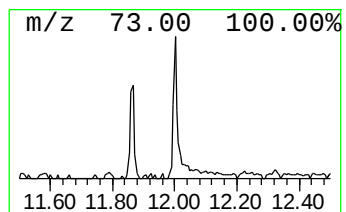
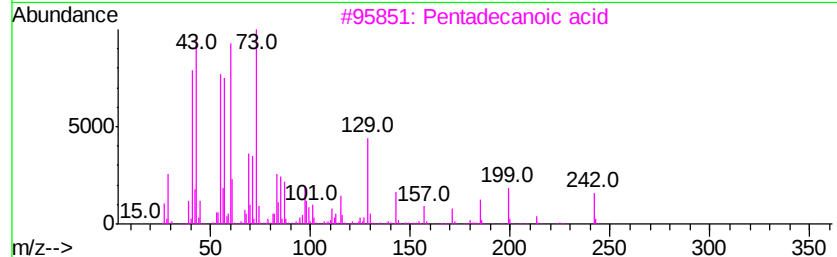
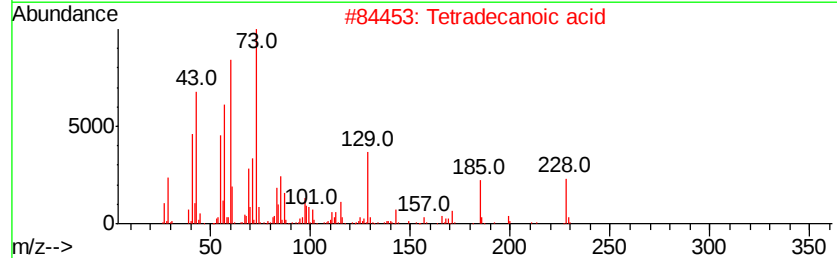
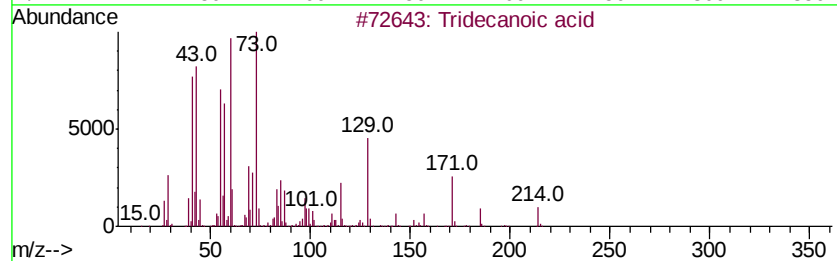
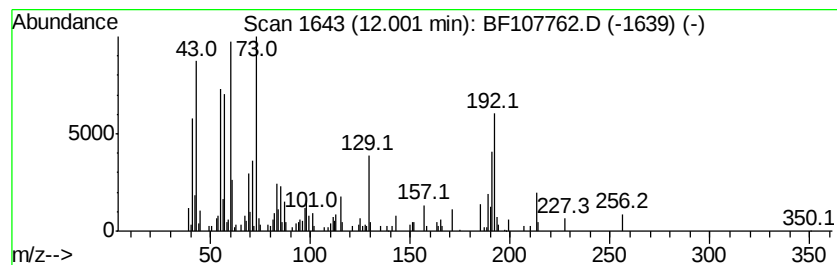
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Tridecanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.00	4.53 ng	188123	Phenanthrene-d10	11.49

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tridecanoic acid	214	C13H26O2	000638-53-9	87
2		Tetradecanoic acid	228	C14H28O2	000544-63-8	47
3		Pentadecanoic acid	242	C15H30O2	001002-84-2	47
4		n-Decanoic acid	172	C10H20O2	000334-48-5	43
5		Dodecanoic acid	200	C12H24O2	000143-07-7	38



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072718\
Data File : BF107762.D
Acq On : 27 Jul 2018 22:37
Operator : JU/SJ
Sample : J4136-02
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CON-SB-01-02-03-0-25

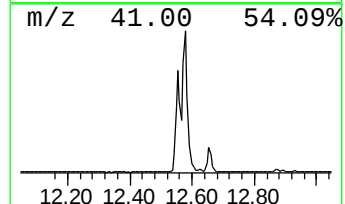
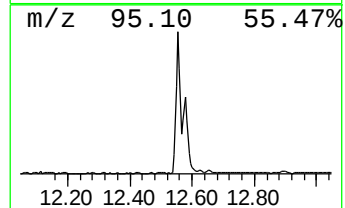
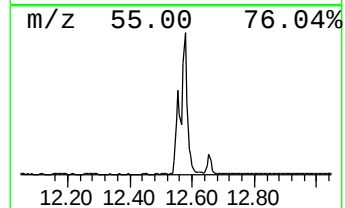
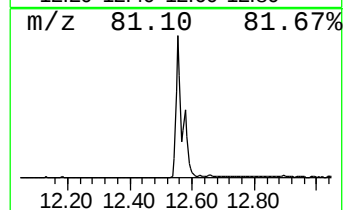
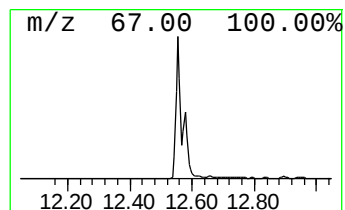
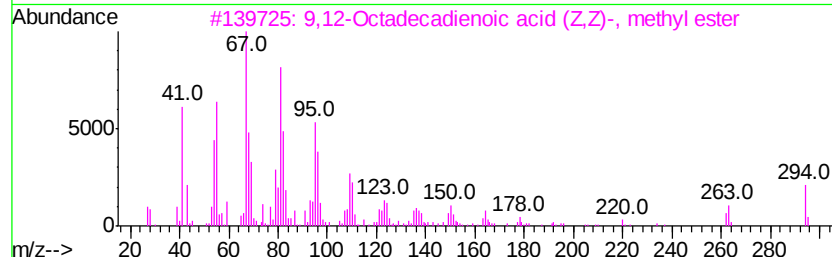
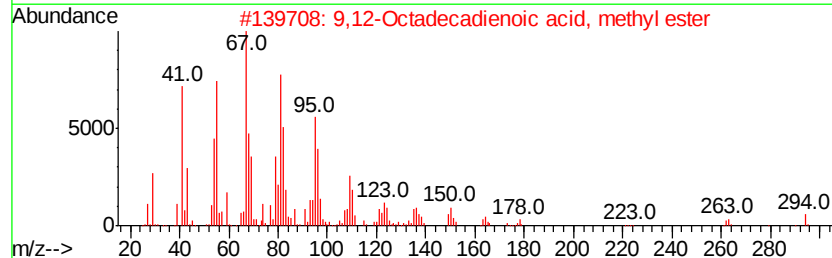
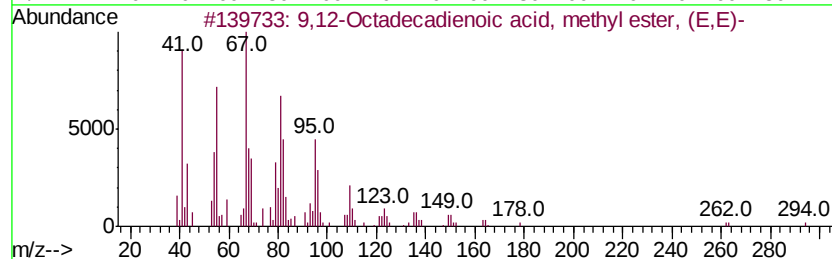
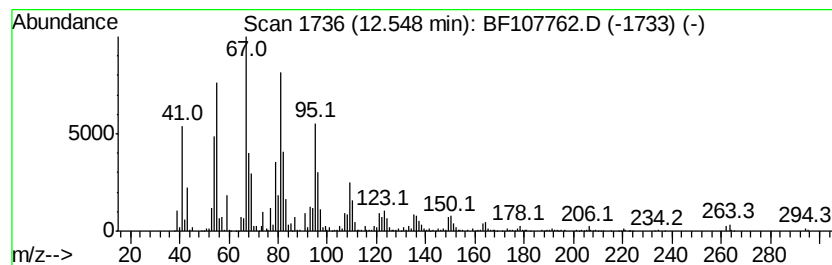
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 9,12-Octadecadienoic acid, ... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.55	96.06 ng	3989080	Phenanthrene-d10	11.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9,12-Octadecadienoic acid, methy...	294	C19H34O2	002566-97-4	99
2		9,12-Octadecadienoic acid, methy...	294	C19H34O2	002462-85-3	99
3		9,12-Octadecadienoic acid (Z,Z)-...	294	C19H34O2	000112-63-0	99
4		10,13-Octadecadienoic acid, meth...	294	C19H34O2	056554-62-2	99
5		8,11-Octadecadienoic acid, methy...	294	C19H34O2	056599-58-7	99



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072718\
Data File : BF107762.D
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Operator : JU/SJ
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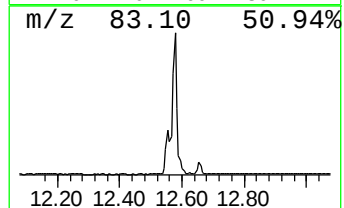
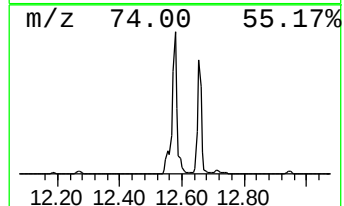
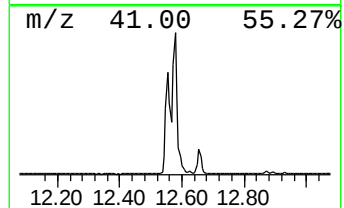
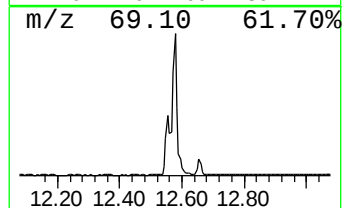
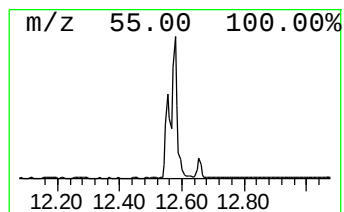
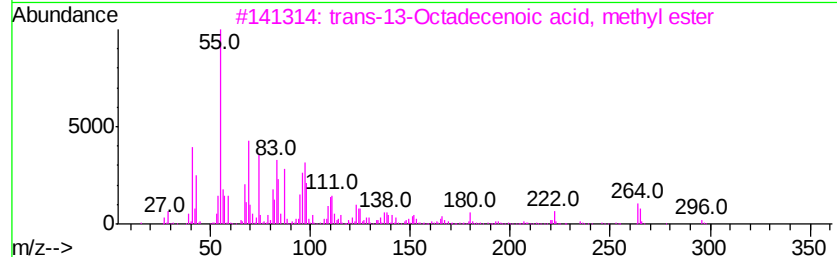
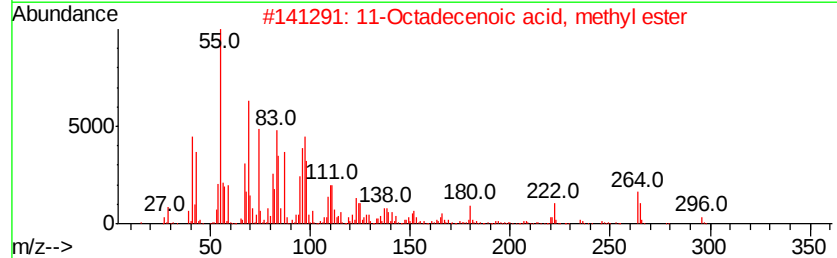
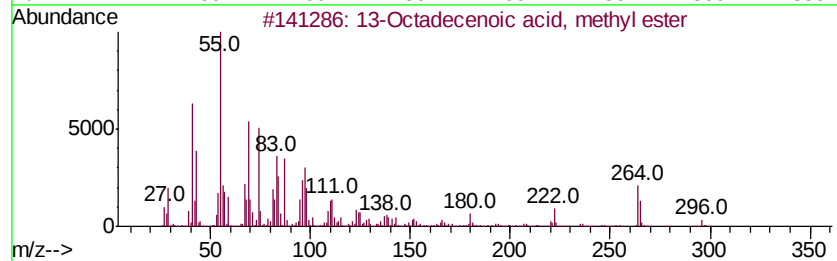
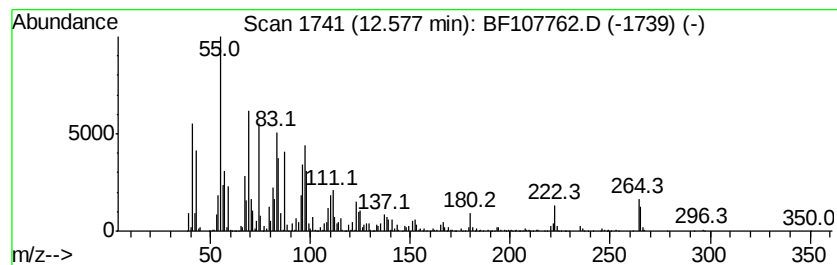
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 13-Octadecenoic acid, methyl... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.58	119.27 ng	4952820	Phenanthrene-d10	11.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	13-Octadecenoic acid, methyl ester	296	C19H36O2	056554-47-3	97
2		11-Octadecenoic acid, methyl ester	296	C19H36O2	052380-33-3	94
3		trans-13-Octadecenoic acid, methyl ester	296	C19H36O2	1000333-61-3	94
4		9-Octadecenoic acid, methyl ester	296	C19H36O2	001937-62-8	94
5		cis-13-Octadecenoic acid, methyl ester	296	C19H36O2	1000333-58-3	94



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072718\
Data File : BF107762.D
Acq On : 27 Jul 2018 22:37
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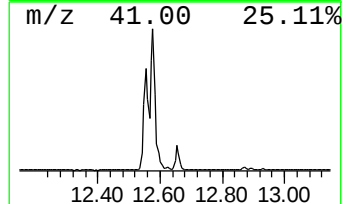
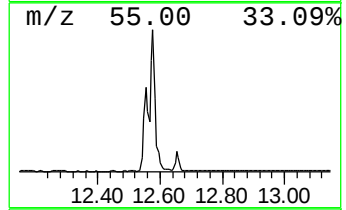
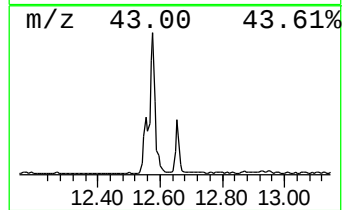
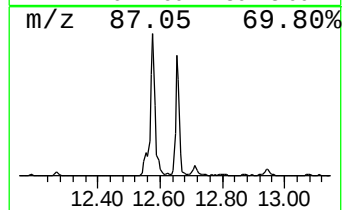
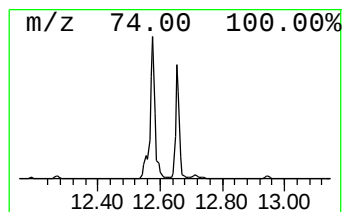
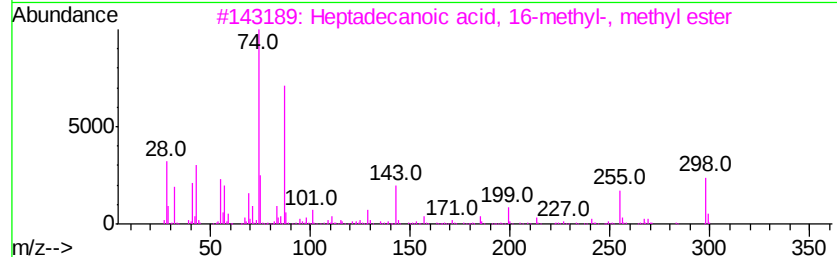
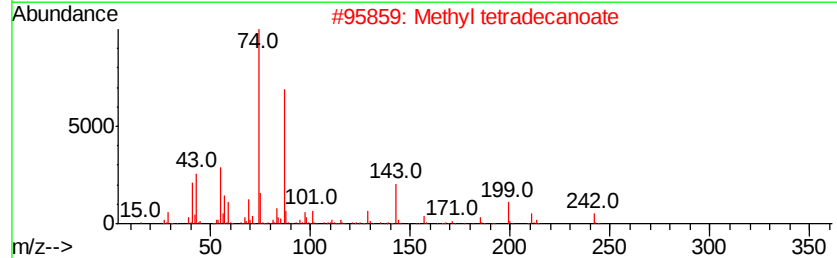
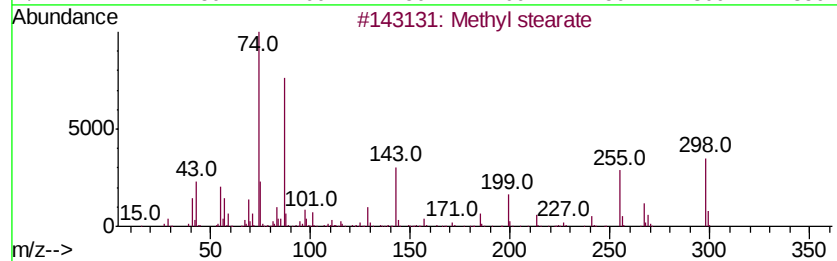
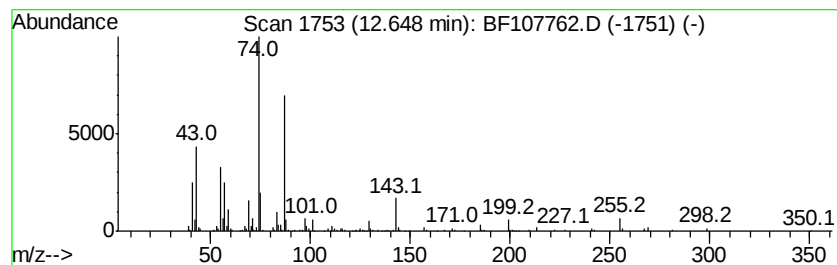
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Methyl stearate Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.65	19.83 ng	823396	Phenanthrene-d10	11.49

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Methyl stearate	298	C19H38O2	000112-61-8	98
2		Methyl tetradecanoate	242	C15H30O2	000124-10-7	94
3		Heptadecanoic acid, 16-methyl-, ...	298	C19H38O2	005129-61-3	91
4		Heptadecanoic acid, methyl ester	284	C18H36O2	001731-92-6	90
5		Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072718\
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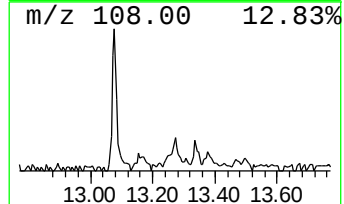
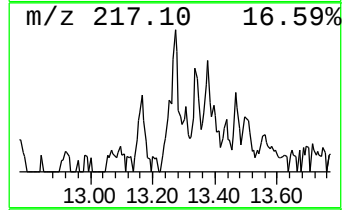
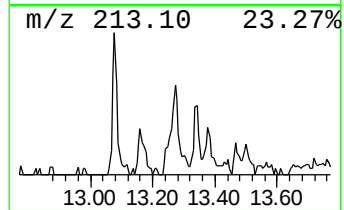
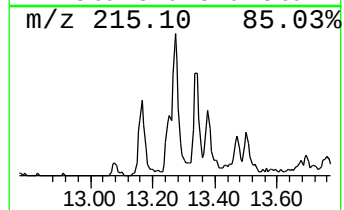
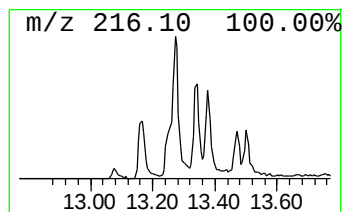
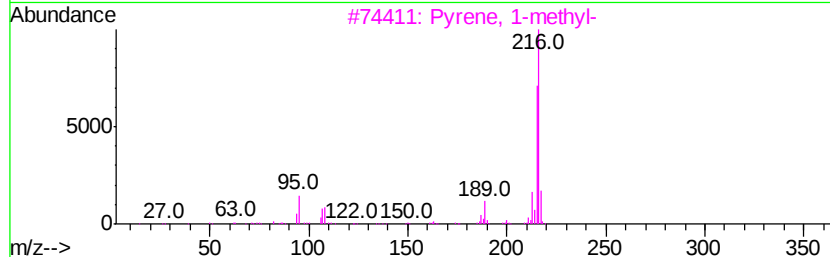
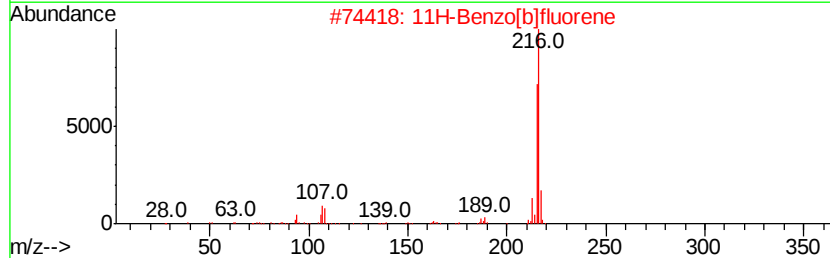
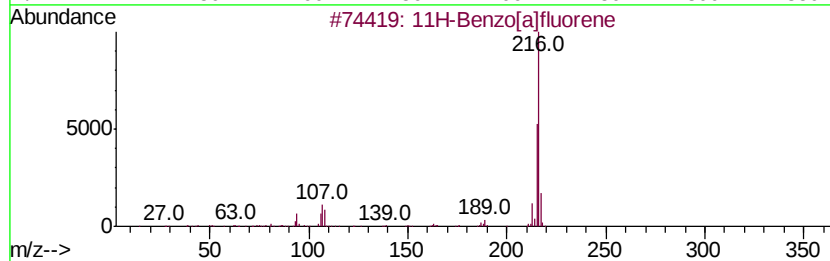
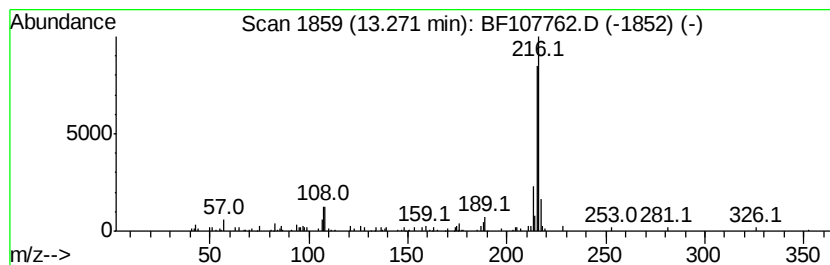
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 11H-Benzo[a]fluorene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.27	2.38 ng	167177	Chrysene-d12	14.15	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	91
2	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	91
3	Pvrene, 1-methyl-	216	C17H12	002381-21-7	90
4	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	83
5	7H-Benzo[c]fluorene	216	C17H12	000205-12-9	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072718\
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Sample : J4136-02
Misc :
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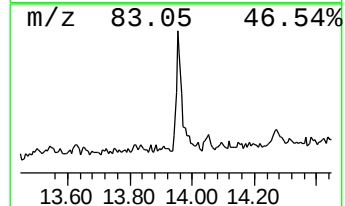
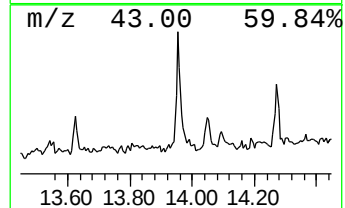
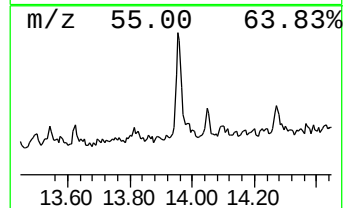
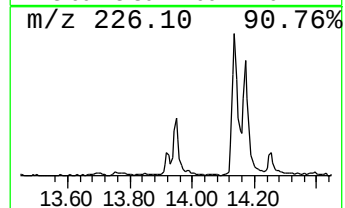
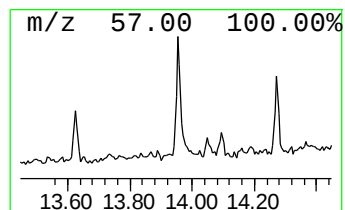
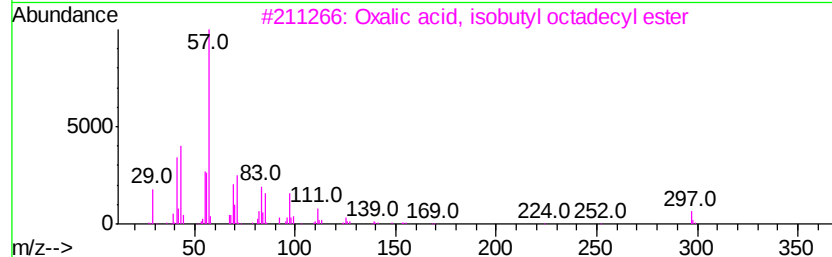
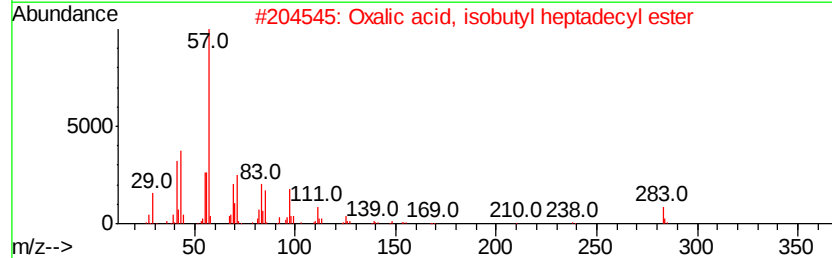
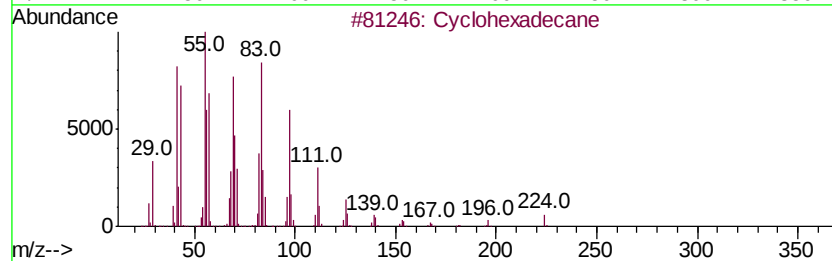
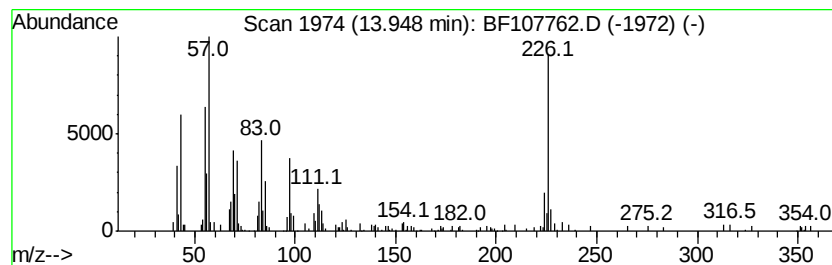
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Cyclohexadecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.		
13.95	4.00 ng	280917	Chrysene-d12	14.15		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cvclohexadecane	224	C16H32	000295-65-8	83
2		Oxalic acid, isobutyl heptadecyl...	384	C23H44O4	1000309-38-2	68
3		Oxalic acid, isobutyl octadecyl ...	398	C24H46O4	1000309-38-3	68
4		Tricosyl trifluoroacetate	436	C25H47F3O2	1000351-75-1	68
5		Dotriacontyl pentafluoropropionate	612	C35H65F5O2	1000351-81-4	68



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072718\
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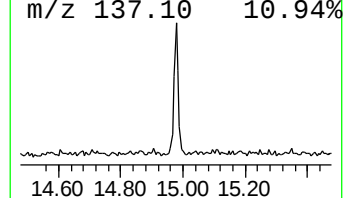
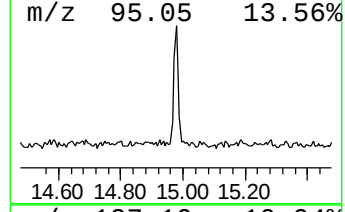
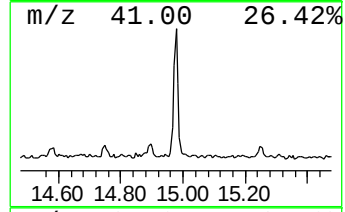
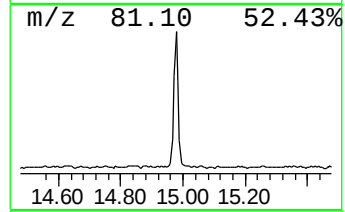
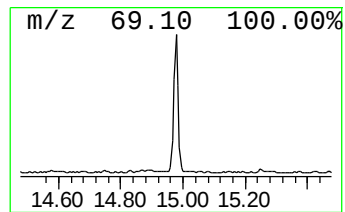
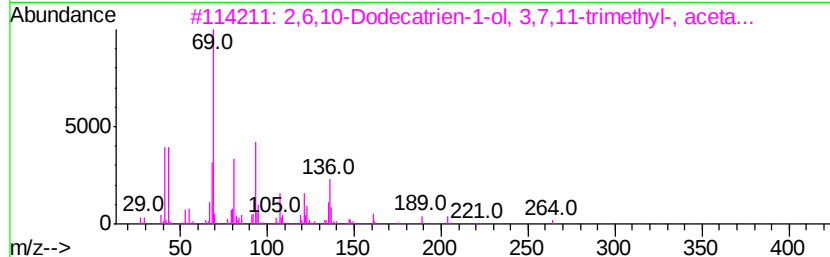
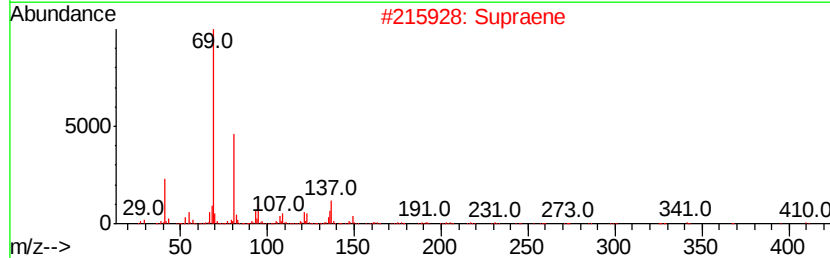
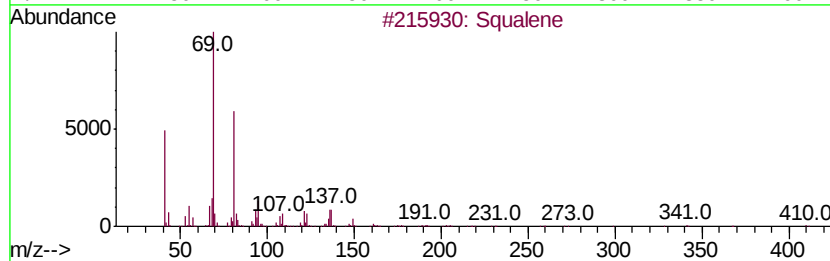
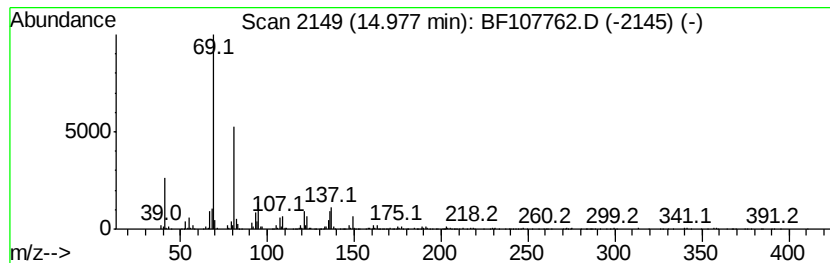
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Squalene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.98	11.53 ng	537605	Perylene-d12	15.68

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Squalene	410	C30H50	000111-02-4	95
2		Supraene	410	C30H50	007683-64-9	91
3		2,6,10-Dodecatrien-1-ol, 3,7,11-...	264	C17H28O2	004128-17-0	64
4		Trifluoroacetyl-lavandulol	250	C12H17F3O2	028673-24-7	64
5		1,5-Heptadiene, 3,3,6-trimethyl-	138	C10H18	035387-63-4	52



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF072718\
Data File : BF107762.D
Acq On : 27 Jul 2018 22:37
Operator : JU/SJ
Sample : J4136-02
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CON-SB-01-02-03-0-25

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF072018.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
2-Pentanone, 4-hy...	5.19	6.7	ng	305742	1	6.96	911876	20.0
unknown6.73	6.73	54.2	ng	2471990	1	6.96	911876	20.0
9-Hexadecenoic ac...	11.79	2.6	ng	106472	4	11.49	830523	20.0
Hexadecanoic acid...	11.87	39.4	ng	1636430	4	11.49	830523	20.0
Tridecanoic acid	12.00	4.5	ng	188123	4	11.49	830523	20.0
9,12-Octadecadien...	12.55	96.1	ng	3989080	4	11.49	830523	20.0
13-Octadecenoic a...	12.58	119.3	ng	4952820	4	11.49	830523	20.0
Methyl stearate	12.65	19.8	ng	823396	4	11.49	830523	20.0
11H-Benzo[a]fluorene	13.27	2.4	ng	167177	5	14.15	1406290	20.0
Cyclohexadecane	13.95	4.0	ng	280917	5	14.15	1406290	20.0
Squalene	14.98	11.5	ng	537605	6	15.68	932300	20.0