

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF030119\
 Data File : BF112799.D
 Acq On : 1 Mar 2019 14:58
 Operator : JU/SJ
 Sample : K1655-02
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 DUP-2

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-BF022719.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	5.357	550	556	559	rBV	3504317	3446898	95.04%	9.852%
2	6.375	724	729	732	rBV	3406789	3606602	99.44%	10.308%
3	6.510	748	752	756	rBV	3703287	3626912	100.00%	10.366%
4	6.745	788	792	795	rBV	1094606	995215	27.44%	2.844%
5	6.898	814	818	822	rBV	3109294	2830505	78.04%	8.090%
6	7.298	881	886	893	rBV	2229826	2254516	62.16%	6.444%
7	8.022	1005	1009	1026	rBV	1308589	1225017	33.78%	3.501%
8	9.098	1188	1192	1202	rBV	3981123	3528766	97.29%	10.086%
9	9.486	1255	1258	1265	rBV	256754	248012	6.84%	0.709%
10	9.769	1302	1306	1315	rVB	1107711	1120437	30.89%	3.202%
11	10.551	1435	1439	1450	rBV	1940417	1976658	54.50%	5.650%
12	10.686	1460	1462	1470	rVB	38914	45411	1.25%	0.130%
13	11.245	1554	1557	1561	rBV	1040609	961616	26.51%	2.748%
14	11.657	1623	1627	1630	rVB	189671	160031	4.41%	0.457%
15	11.786	1645	1649	1659	rBV	730877	689701	19.02%	1.971%
16	11.963	1676	1679	1688	rVB	37071	51506	1.42%	0.147%
17	12.351	1743	1745	1748	rVB2	47616	49067	1.35%	0.140%
18	12.445	1758	1761	1766	rBV2	122509	122581	3.38%	0.350%
19	12.486	1766	1768	1777	rVB6	32257	48774	1.34%	0.139%
20	12.568	1778	1782	1792	rBV	303837	352140	9.71%	1.006%
21	12.663	1795	1798	1803	rBV2	158171	167041	4.61%	0.477%
22	12.710	1803	1806	1810	rVB2	52118	74072	2.04%	0.212%
23	12.827	1822	1826	1830	rBV	2889321	2667516	73.55%	7.624%
24	13.074	1865	1868	1872	rVB	43681	42401	1.17%	0.121%
25	13.368	1914	1918	1921	rBV	238726	200205	5.52%	0.572%
26	13.415	1923	1926	1929	rVV	61742	51142	1.41%	0.146%
27	13.733	1976	1980	1989	rVB	464347	556599	15.35%	1.591%
28	13.874	1999	2004	2007	rBV	1179031	1123828	30.99%	3.212%
29	14.057	2032	2035	2038	rBV	132314	136801	3.77%	0.391%
30	14.362	2084	2087	2092	rBV	202339	192179	5.30%	0.549%
31	14.656	2134	2137	2144	rVB	191486	216027	5.96%	0.617%
32	14.727	2146	2149	2153	rVB	108568	103721	2.86%	0.296%
33	14.980	2189	2192	2200	rVB	195208	209443	5.77%	0.599%
34	15.280	2239	2243	2248	rVB	927789	1084027	29.89%	3.098%

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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

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35	15.339	2249	2253	2258	rVB	161738	207037	5.71%	0.592%
36	15.745	2318	2322	2326	rBV	145301	225893	6.23%	0.646%
37	15.786	2326	2329	2339	rVB2	128236	222154	6.13%	0.635%
38	16.215	2398	2402	2408	rVB	112044	167701	4.62%	0.479%

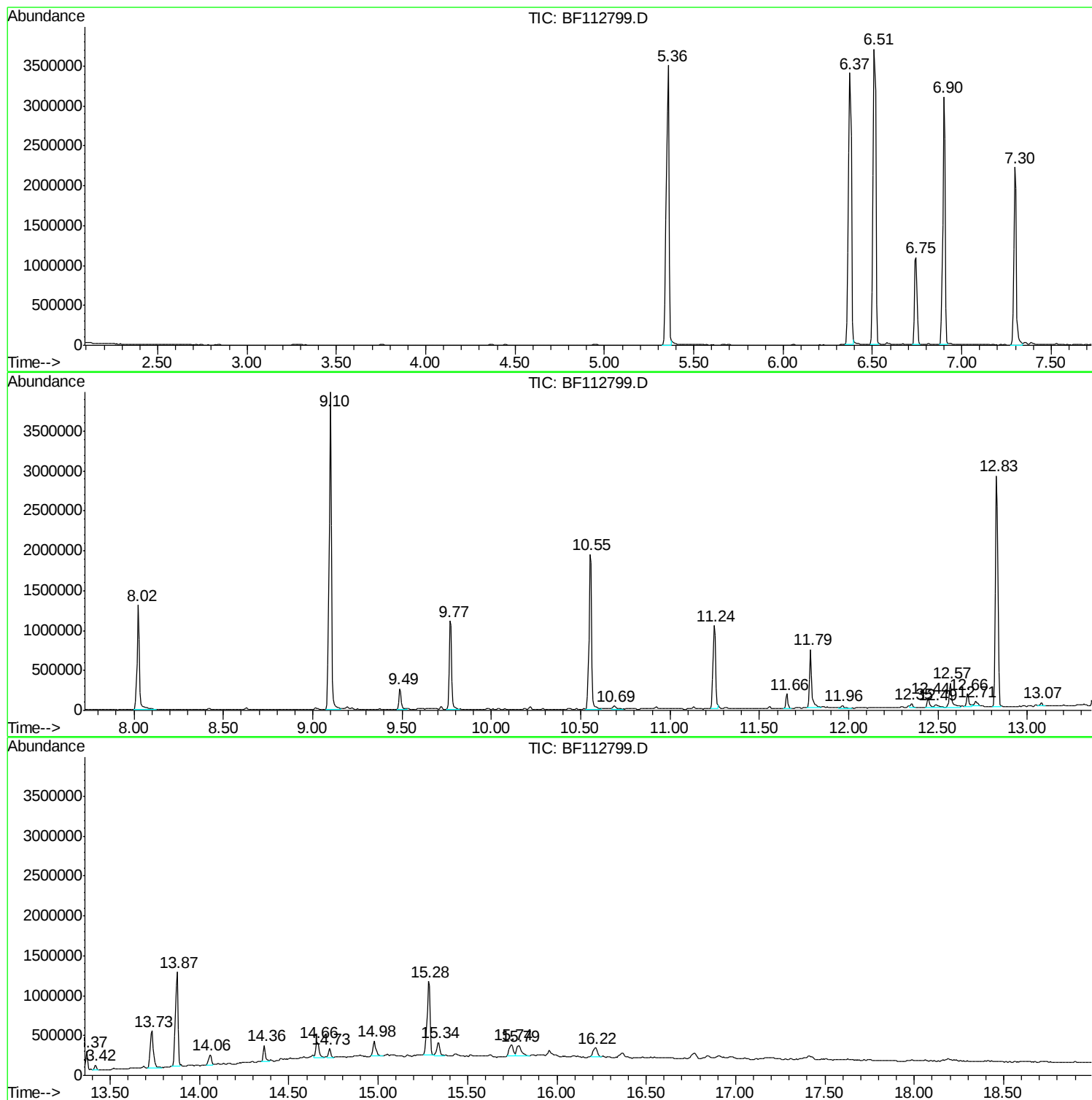
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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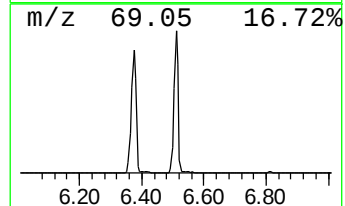
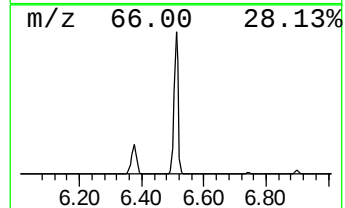
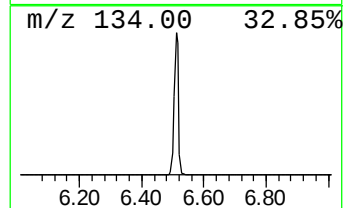
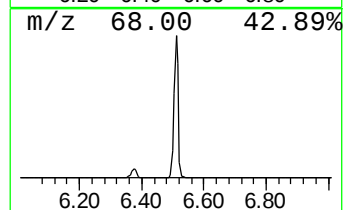
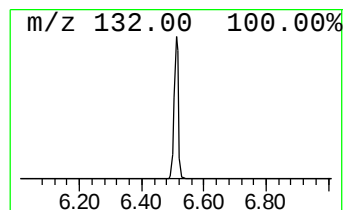
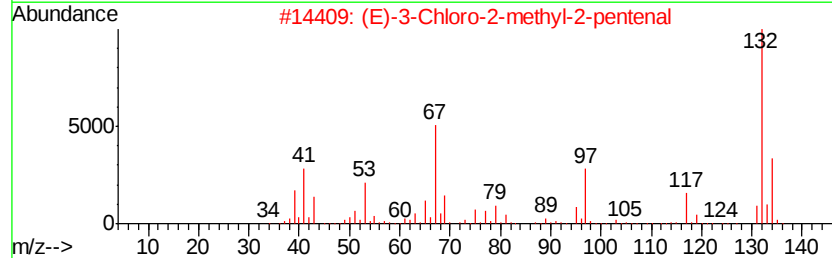
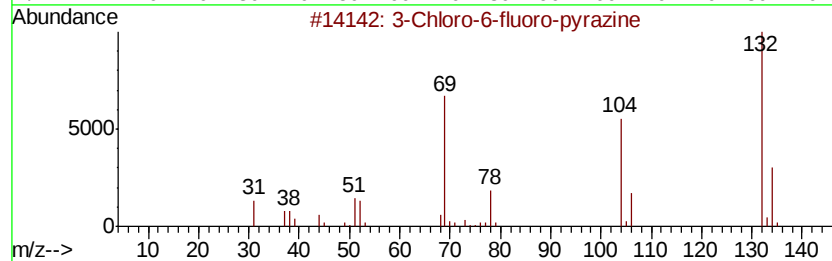
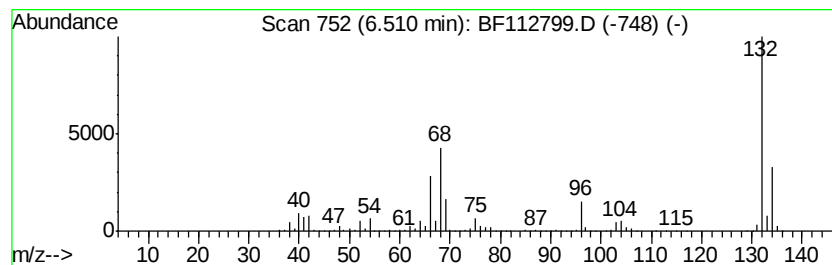
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 1 unknown6.51 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.51	72.89 ng	3626910	1,4-Dichlorobenzene-d4	6.75

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	16
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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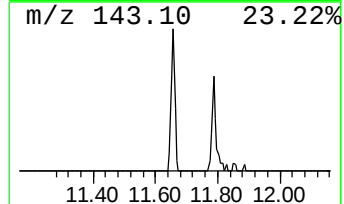
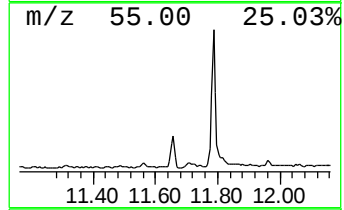
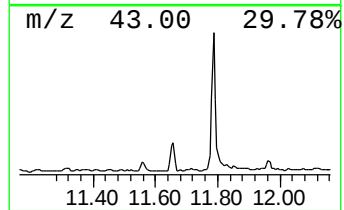
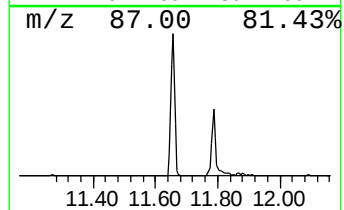
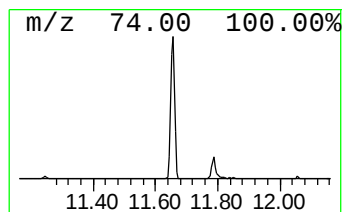
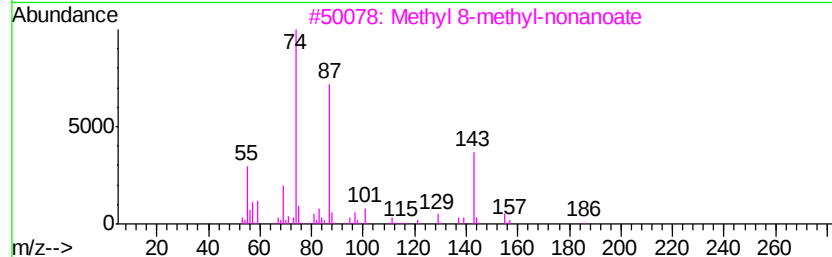
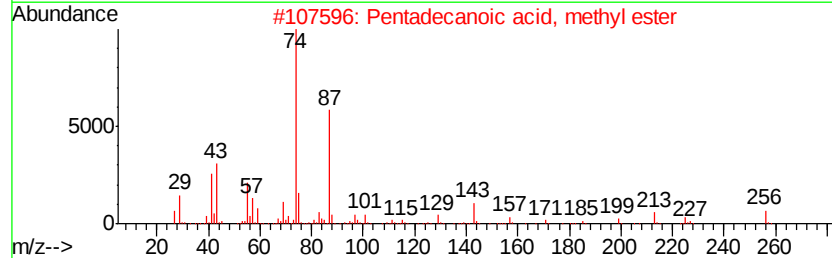
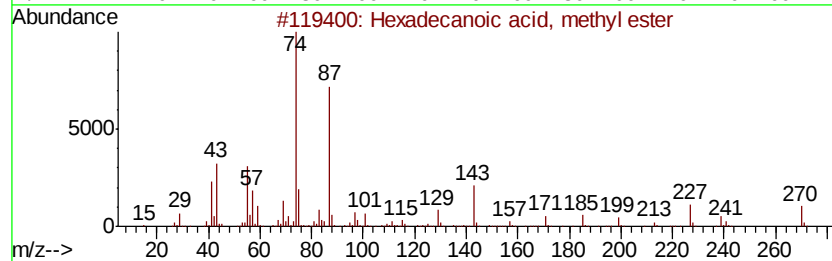
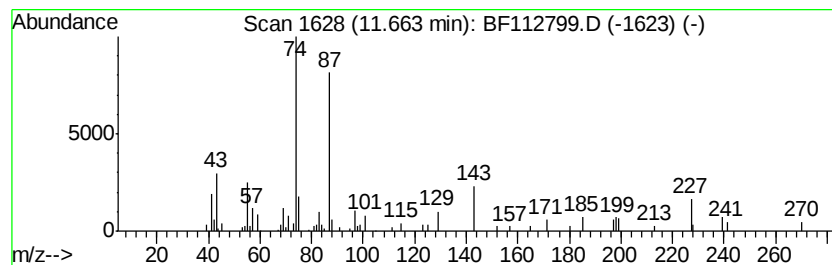
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Hexadecanoic acid, methyl e... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.66	3.33 ng	160031	Phenanthrene-d10	11.24

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	91
2			Pentadecanoic acid, methyl ester	256	C16H32O2	007132-64-1	86
3			Methyl 8-methyl-nonanoate	186	C11H22O2	1000336-43-6	81
4			Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	80
5			Methyl 11-methyl-dodecanoate	228	C14H28O2	1000336-45-1	80



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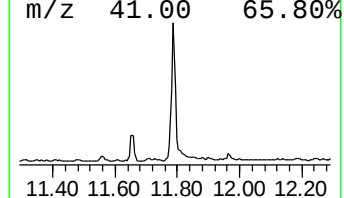
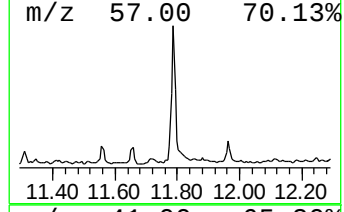
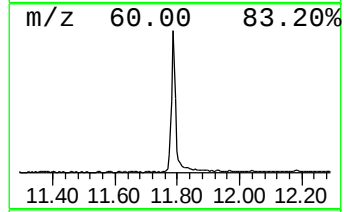
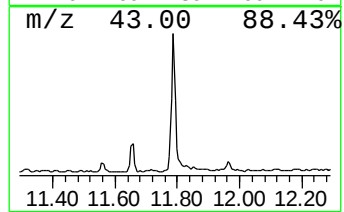
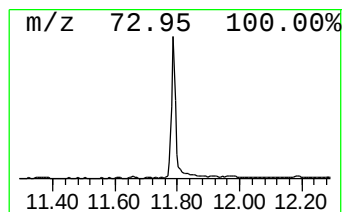
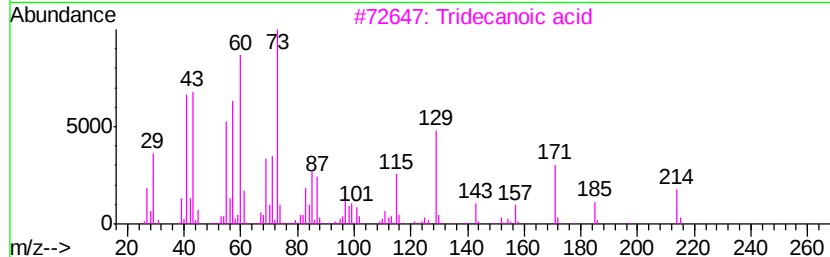
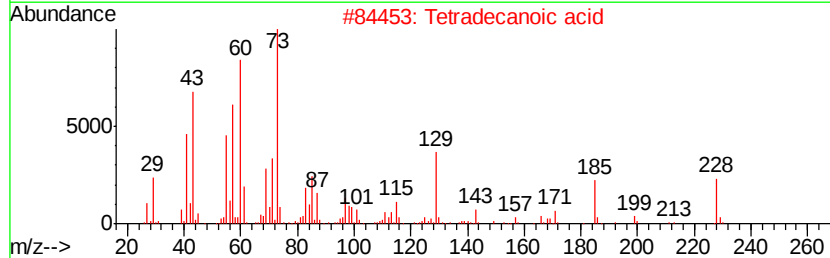
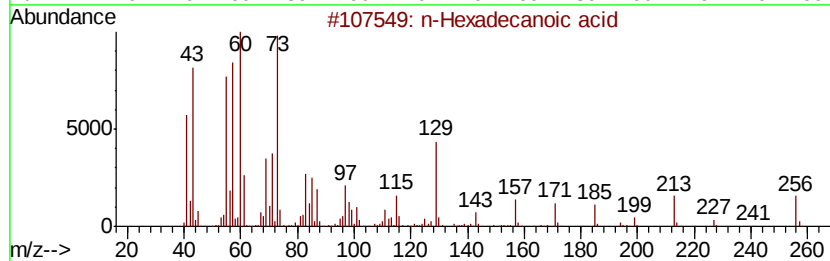
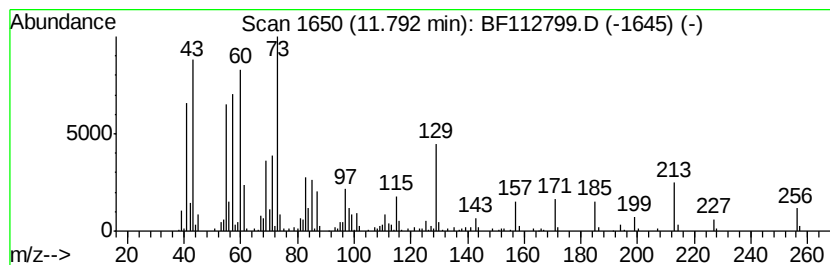
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 4 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.79	14.34 ng	689701	Phenanthrene-d10	11.24

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		Tetradecanoic acid	228	C14H28O2	000544-63-8	91
3		Tridecanoic acid	214	C13H26O2	000638-53-9	90
4		Pentadecanoic acid	242	C15H30O2	001002-84-2	90
5		Dodecanoic acid	200	C12H24O2	000143-07-7	64



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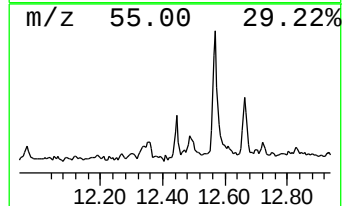
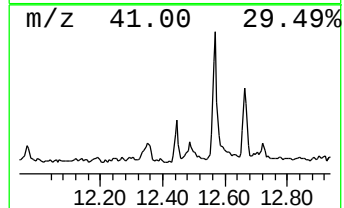
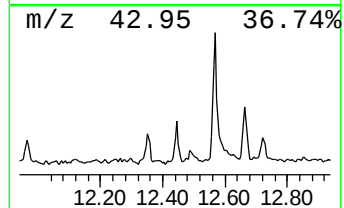
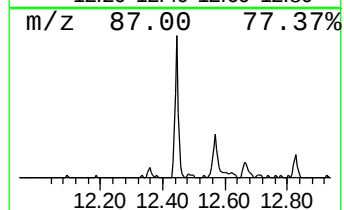
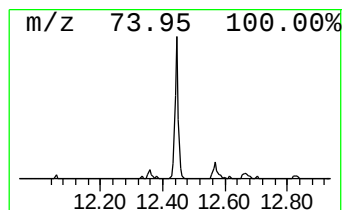
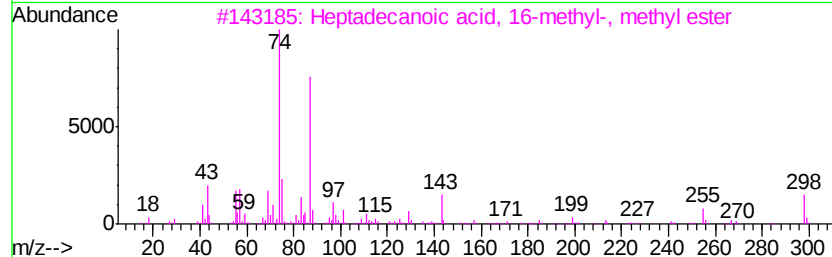
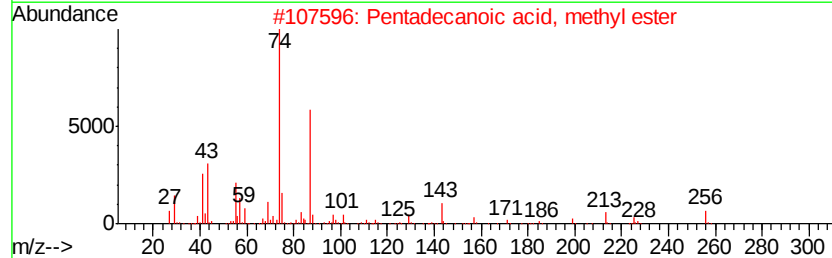
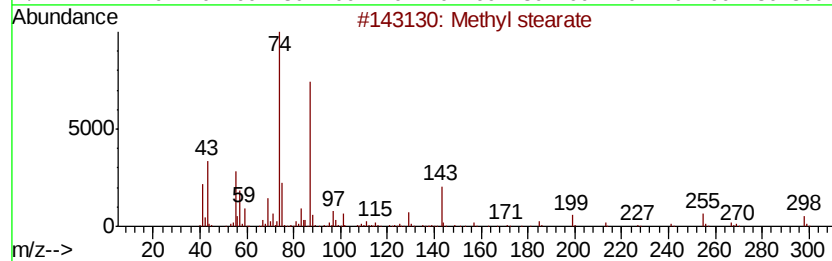
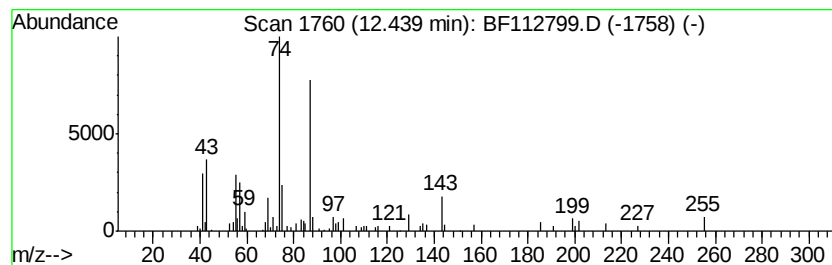
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Methyl stearate Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.44	2.55 ng	122581	Phenanthrene-d10	11.24
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Methyl stearate	298 C19H38O2	000112-61-8 96
2		Pentadecanoic acid, methyl ester	256 C16H32O2	007132-64-1 81
3		Heptadecanoic acid, 16-methyl-, ...	298 C19H38O2	005129-61-3 80
4		Methyl 13-methyltetradecanoate	256 C16H32O2	1000336-31-4 72
5		Undecanoic acid, methyl ester	200 C12H24O2	001731-86-8 72



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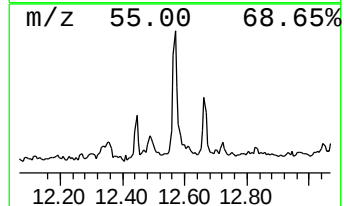
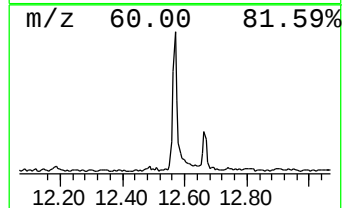
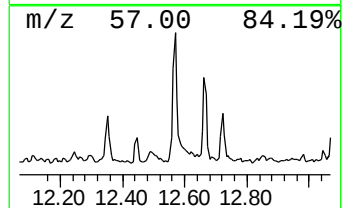
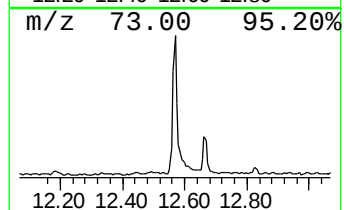
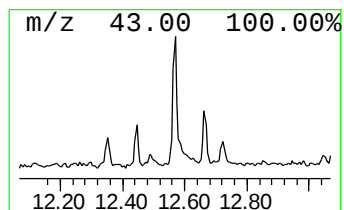
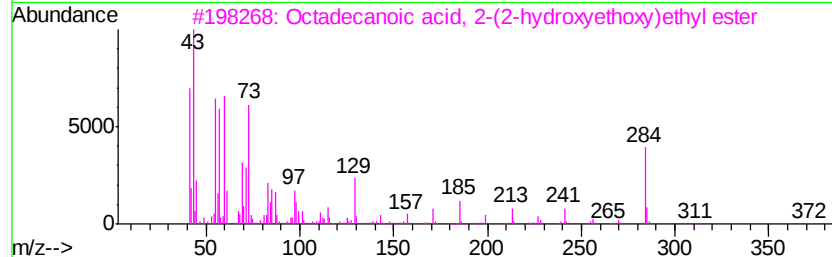
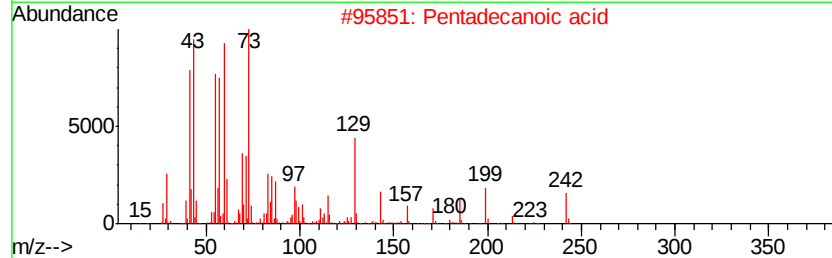
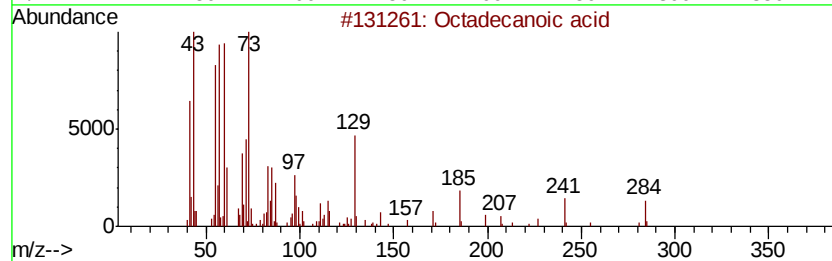
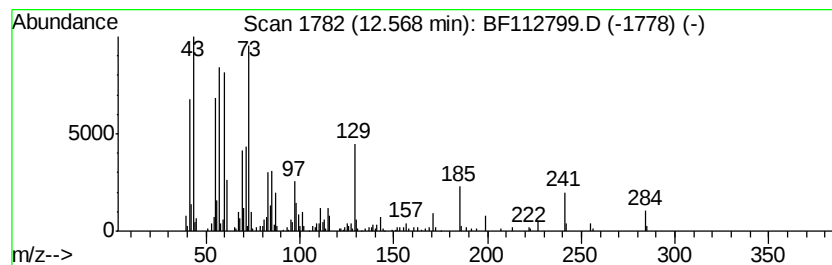
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 6 Octadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.57	6.27 ng	352140	Chrysene-d12	13.87

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	91
3		Octadecanoic acid, 2-(2-hydroxyve...	372	C22H44O4	000106-11-6	91
4		Tetradecanoic acid	228	C14H28O2	000544-63-8	83
5		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	62



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF030119\
Data File : BF112799.D
Acq On : 1 Mar 2019 14:58
Operator : JU/SJ
Sample : K1655-02
Misc :
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Instrument :
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ClientSampleId :
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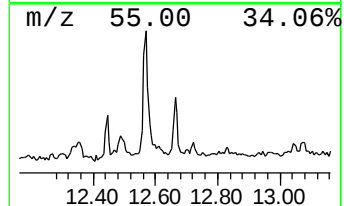
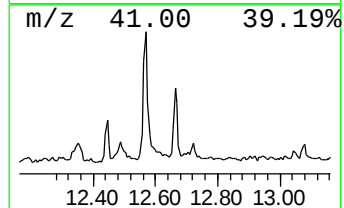
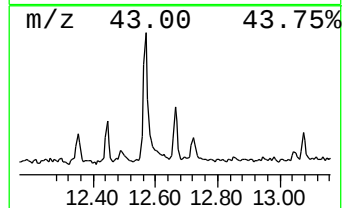
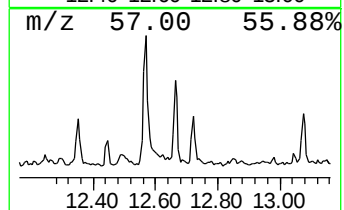
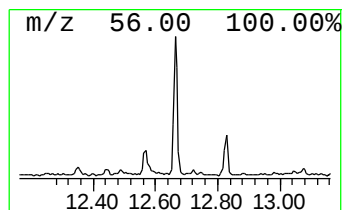
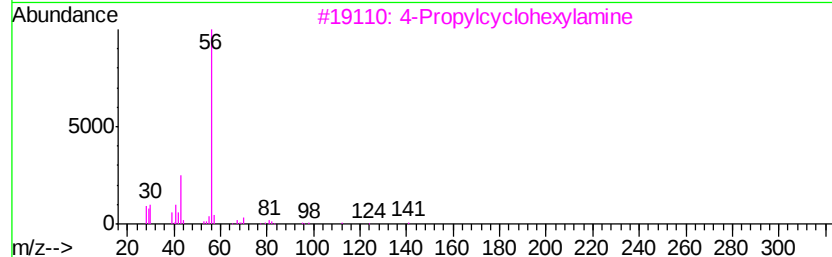
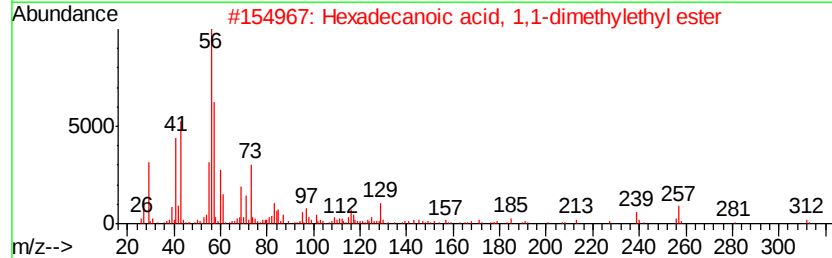
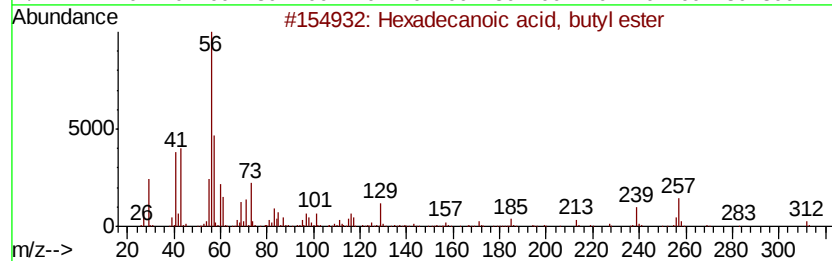
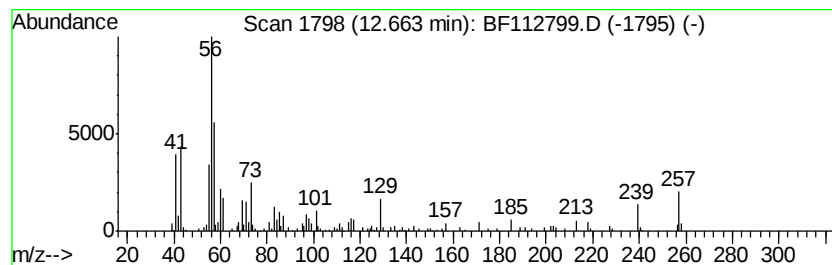
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Hexadecanoic acid, butyl ester Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.66	2.97 ng	167041	Chrysene-d12	13.87

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecanoic acid, butyl ester	312	C20H40O2	000111-06-8	91
2		Hexadecanoic acid, 1,1-dimethyle...	312	C20H40O2	031158-91-5	72
3		4-Propylcyclohexylamine	141	C9H19N	102653-37-2	27
4		1-Hexene, 2,5-dimethyl-	112	C8H16	006975-92-4	27
5		1-Pentene, 2,4-dimethyl-	98	C7H14	002213-32-3	27



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF030119\
Data File : BF112799.D
Acq On : 1 Mar 2019 14:58
Operator : JU/SJ
Sample : K1655-02
Misc :
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ClientSampleId :
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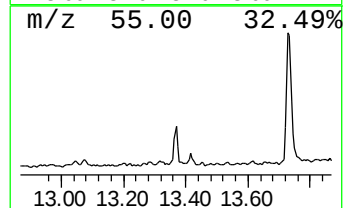
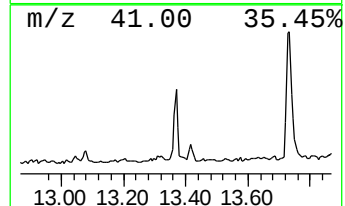
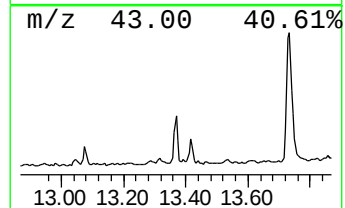
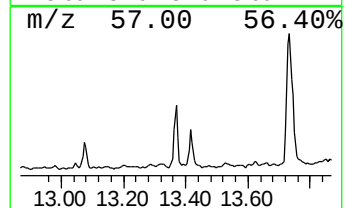
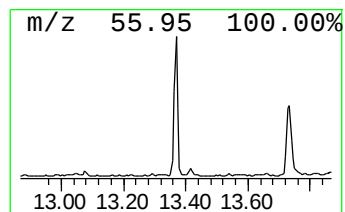
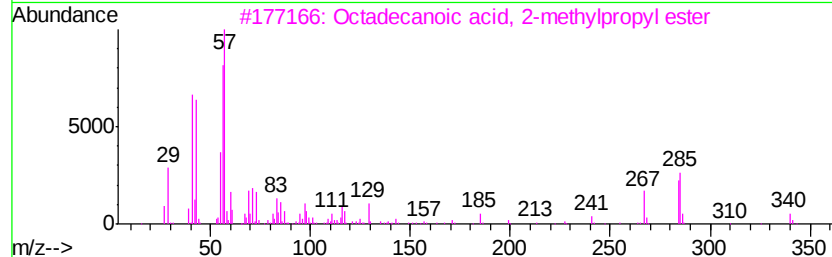
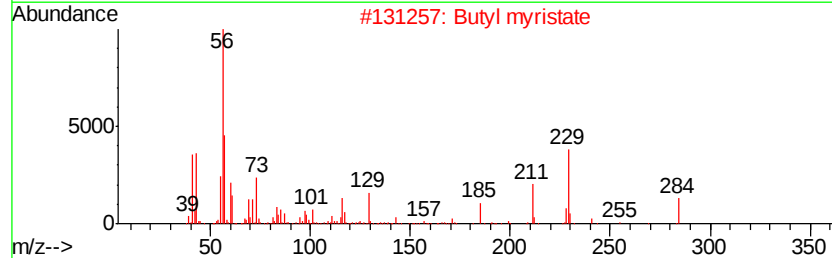
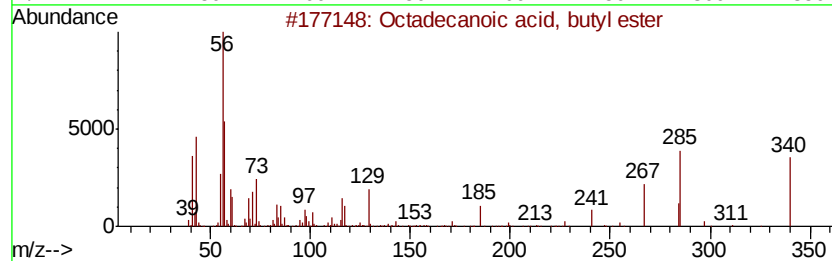
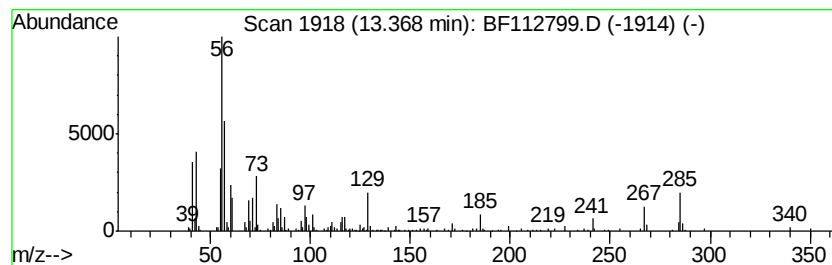
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Octadecanoic acid, butyl ester Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.37	3.56 ng	200205	Chrysene-d12	13.87

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecanoic acid, butyl ester	340	C22H44O2	000123-95-5	95
2		Butyl myristate	284	C18H36O2	000110-36-1	72
3		Octadecanoic acid, 2-methylpropyl...	340	C22H44O2	000646-13-9	46
4		Decane, 2,2,3-trimethyl-	184	C13H28	062338-09-4	35
5		Cyclobutane, 1-hexyl-2,3-dimethyl-	168	C12H24	055170-84-8	35



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF030119\
Data File : BF112799.D
Acq On : 1 Mar 2019 14:58
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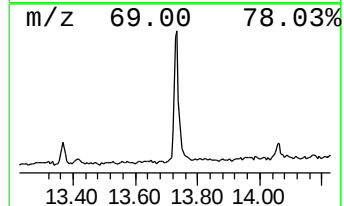
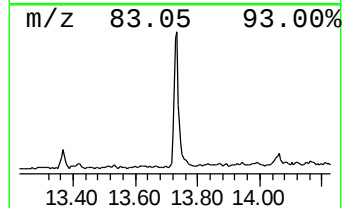
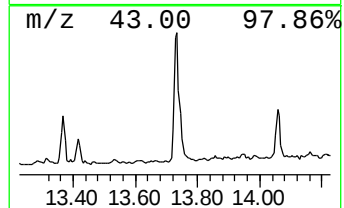
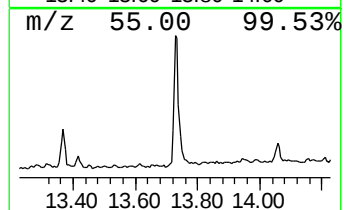
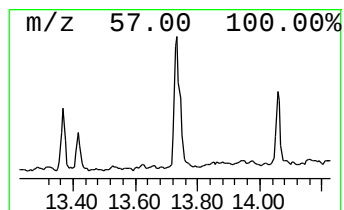
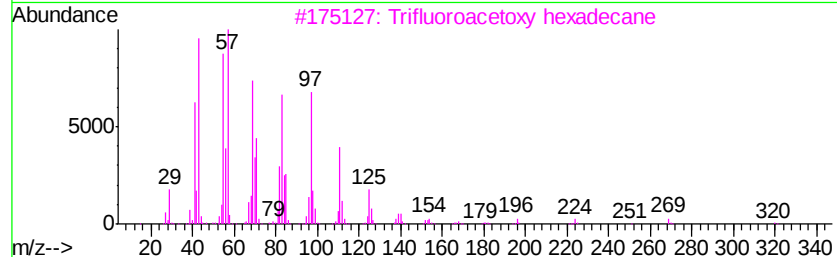
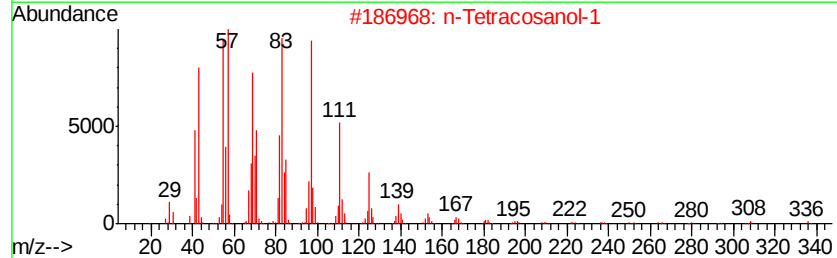
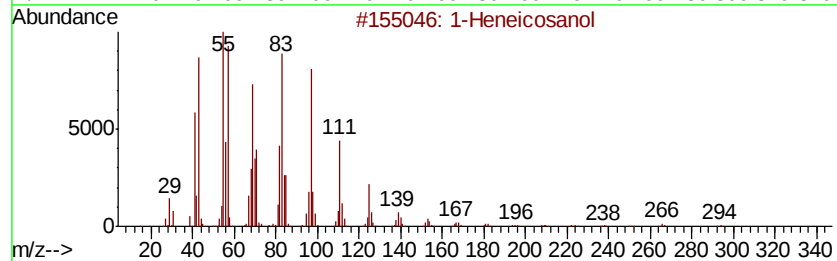
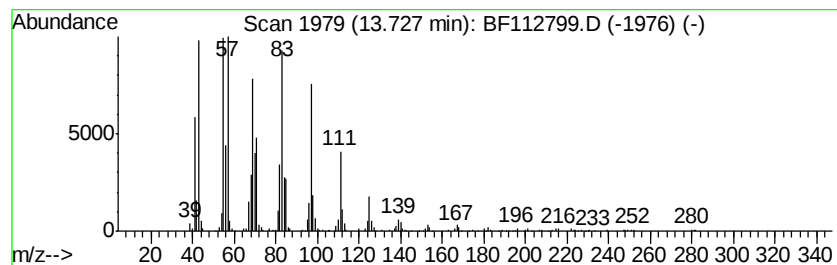
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 1-Heneicosanol Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.73	9.91 ng	556599	Chrysene-d12	13.87

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Heneicosanol	312	C21H44O	015594-90-8	95
2		n-Tetracosanol-1	354	C24H50O	000506-51-4	94
3		Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	94
4		1-Heneicosyl formate	340	C22H44O2	077899-03-7	93
5		n-Nonadecanol-1	284	C19H40O	001454-84-8	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF030119\
Data File : BF112799.D
Acq On : 1 Mar 2019 14:58
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ClientSampleId :
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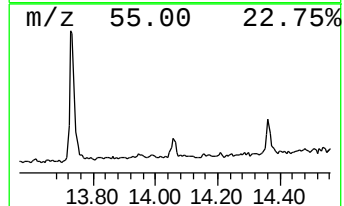
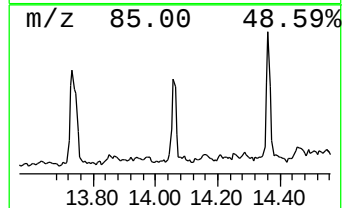
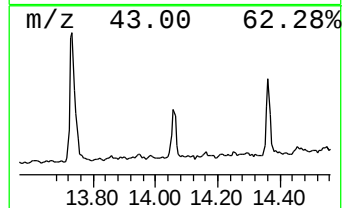
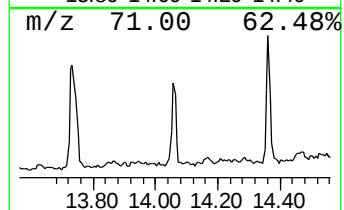
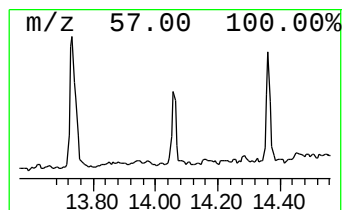
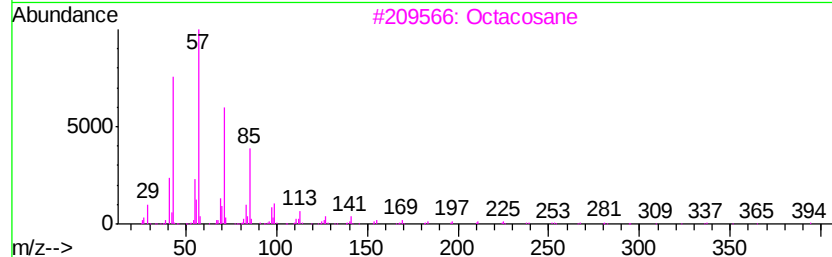
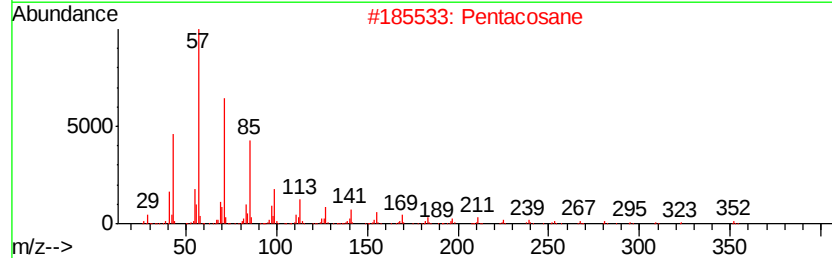
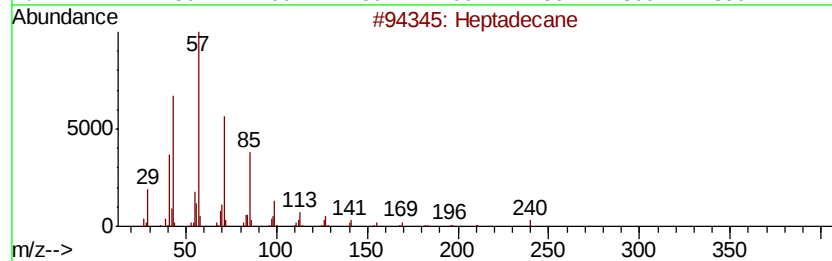
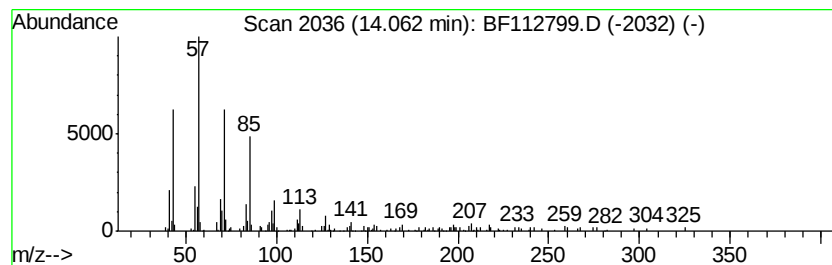
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Heptadecane Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.06	2.43 ng	136801	Chrysene-d12	13.87

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane	240	C17H36	000629-78-7	91
2		Pentacosane	352	C25H52	000629-99-2	87
3		Octacosane	394	C28H58	000630-02-4	87
4		Heptacosane	380	C27H56	000593-49-7	87
5		Hexacosane	366	C26H54	000630-01-3	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF030119\
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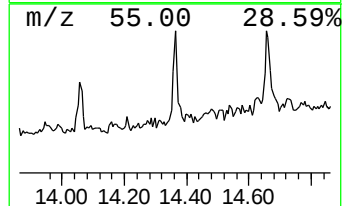
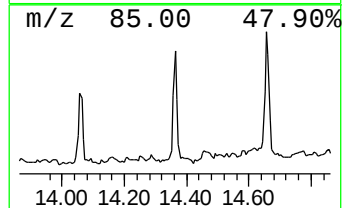
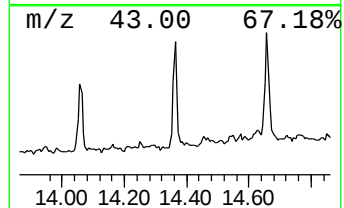
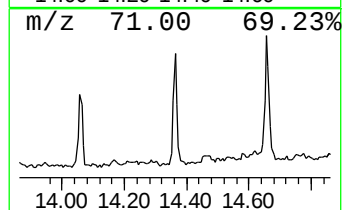
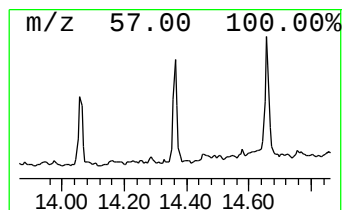
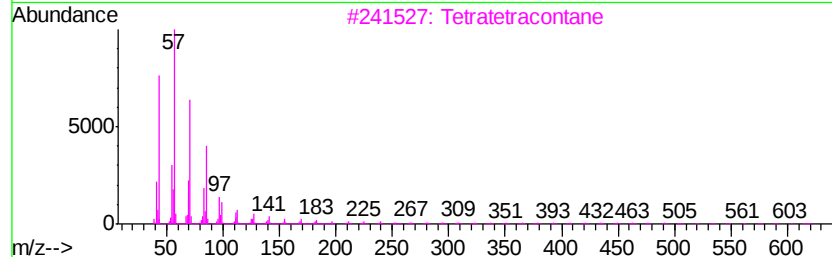
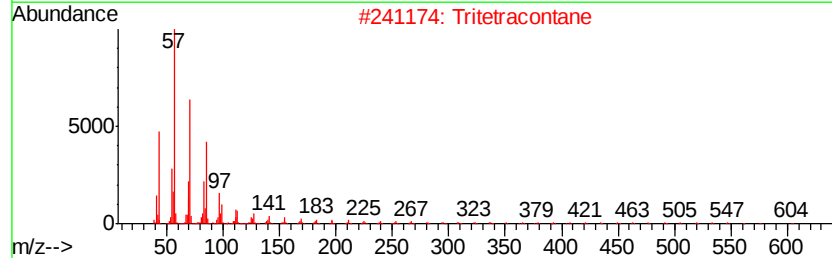
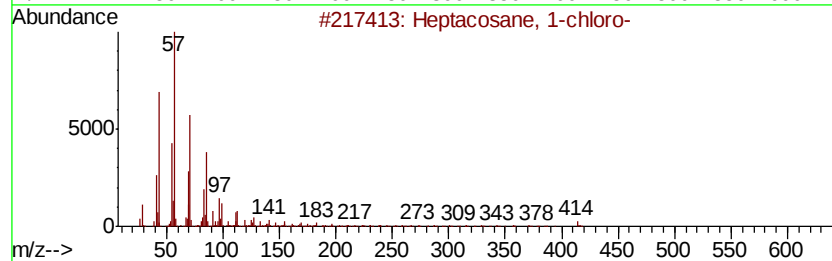
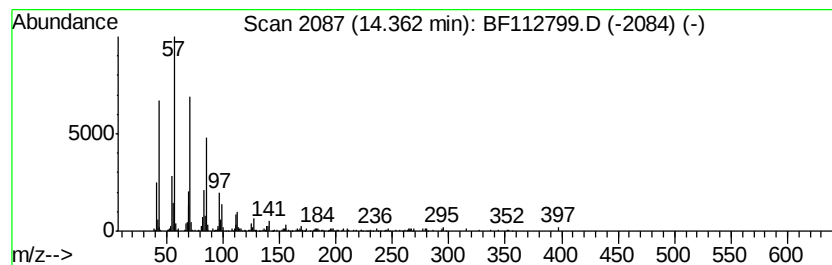
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 Heptacosane, 1-chloro- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.36	3.42 ng	192179	Chrysene-d12	13.87

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptacosane, 1-chloro-	414	C27H55Cl	062016-79-9	91
2			Tritetracontane	605	C43H88	007098-21-7	91
3			Tetratetracontane	619	C44H90	007098-22-8	87
4			Sulfurous acid, dodecyl 2-propyl...	292	C15H32O3S	1000309-12-3	87
5			Tetradecane	198	C14H30	000629-59-4	86



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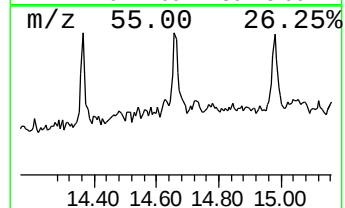
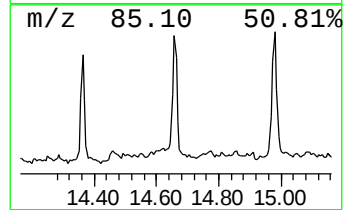
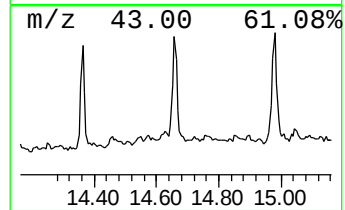
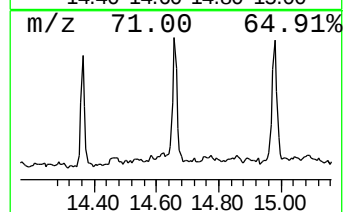
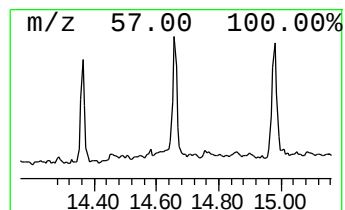
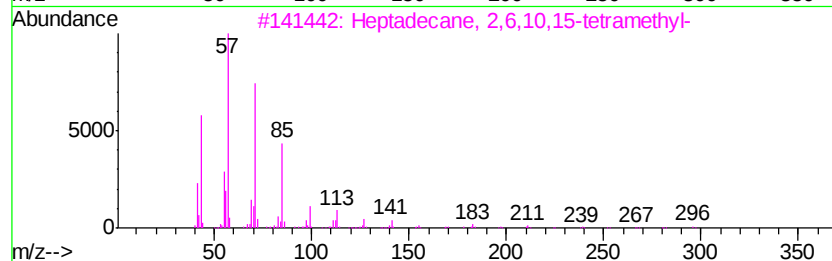
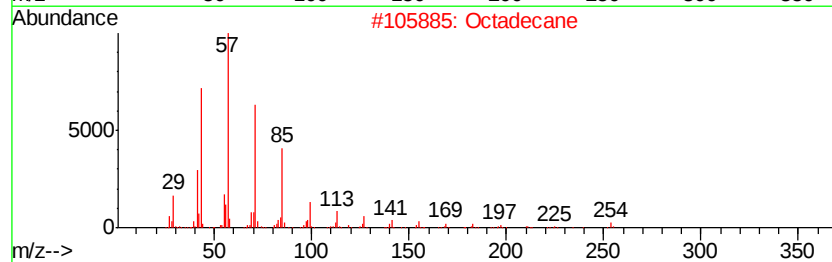
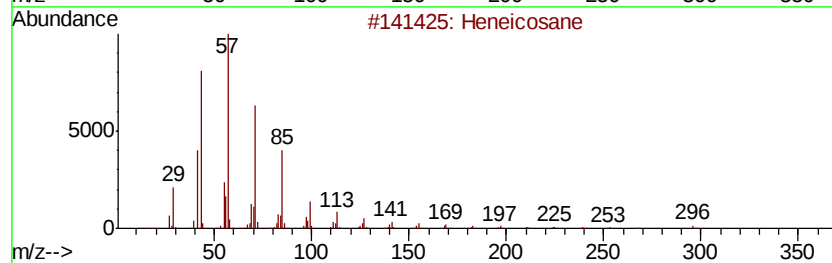
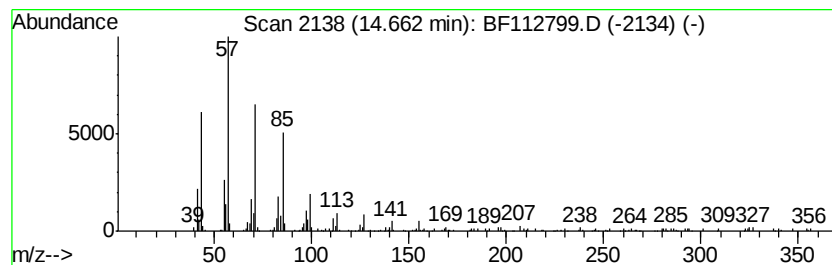
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 Heneicosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.66	3.99 ng	216027	Perylene-d12	15.28

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	93
2		Octadecane	254	C18H38	000593-45-3	93
3		Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	87
4		Octacosane	394	C28H58	000630-02-4	86
5		2-methyloctacosane	408	C29H60	1000376-72-8	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF030119\
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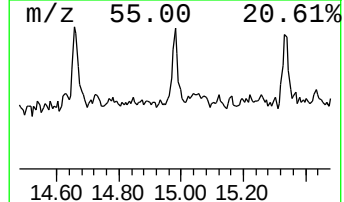
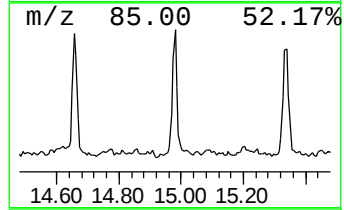
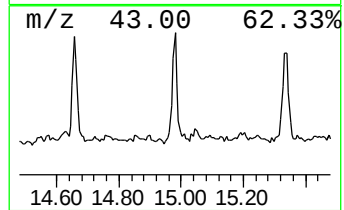
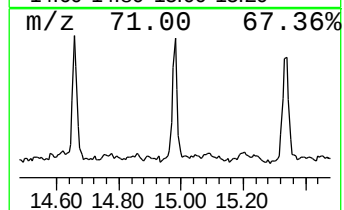
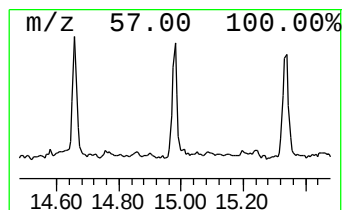
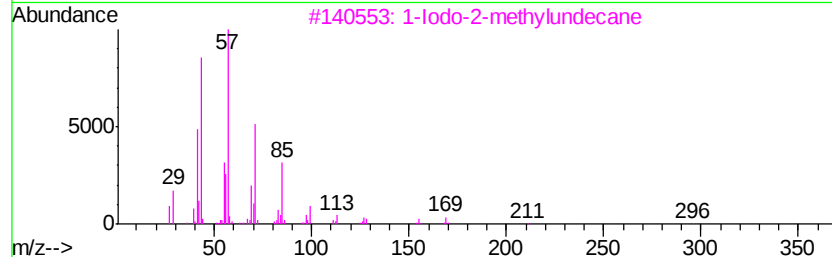
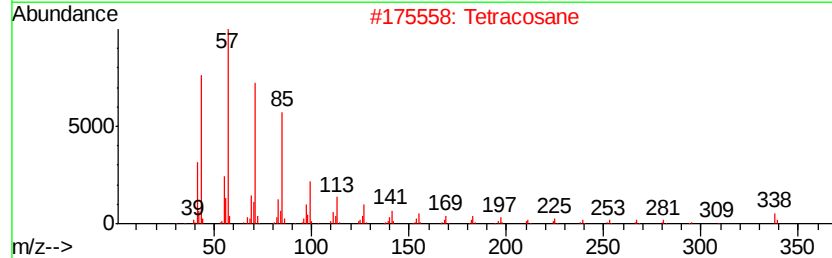
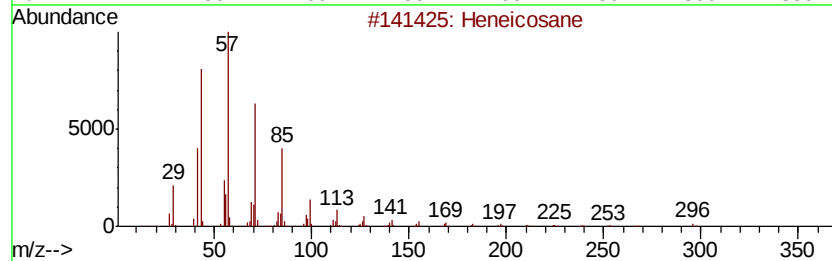
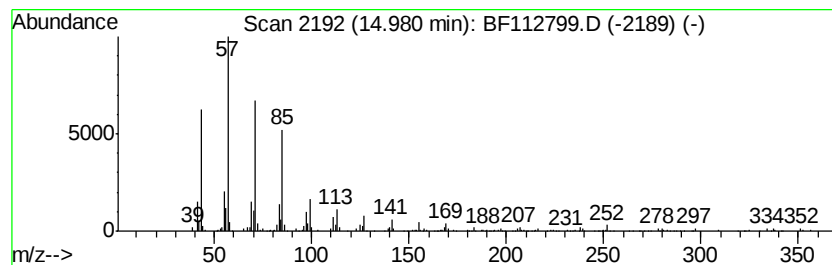
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 1-Iodo-2-methylundecane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.98	3.86 ng	209443	Perylene-d12	15.28

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	97
2			Tetracosane	338	C24H50	000646-31-1	95
3			1-Iodo-2-methylundecane	296	C12H25I	073105-67-6	93
4			Eicosane, 10-methyl-	296	C21H44	054833-23-7	93
5			Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF030119\
 Data File : BF112799.D
 Acq On : 1 Mar 2019 14:58
 Operator : JU/SJ
 Sample : K1655-02
 Misc :
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Instrument :
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 ClientSampleId :
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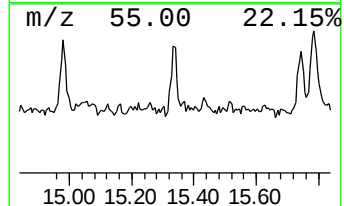
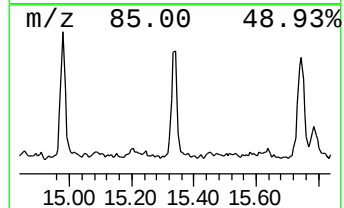
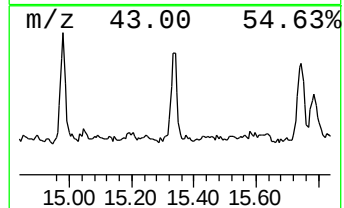
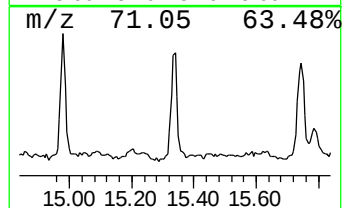
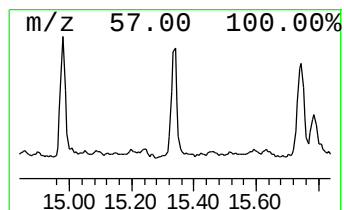
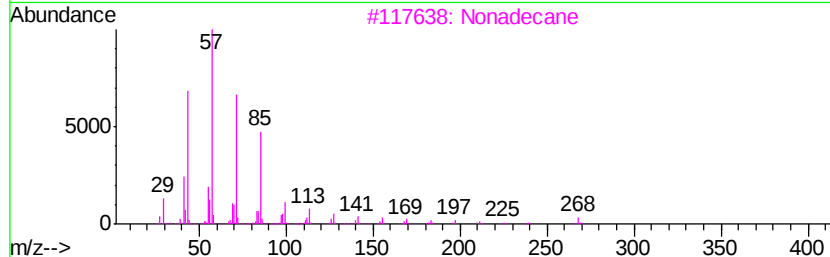
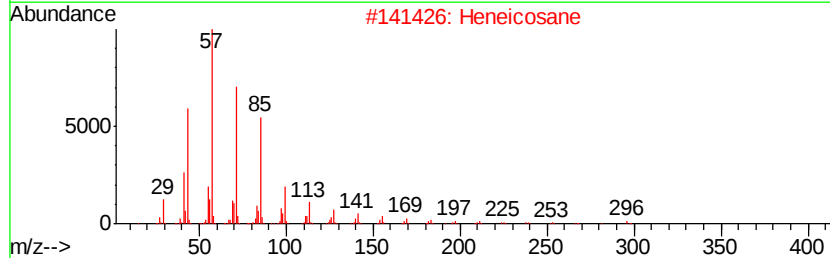
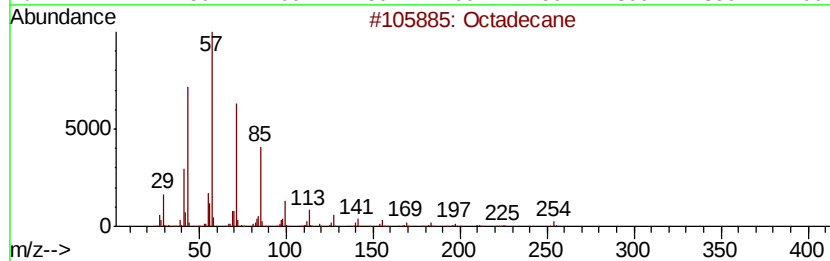
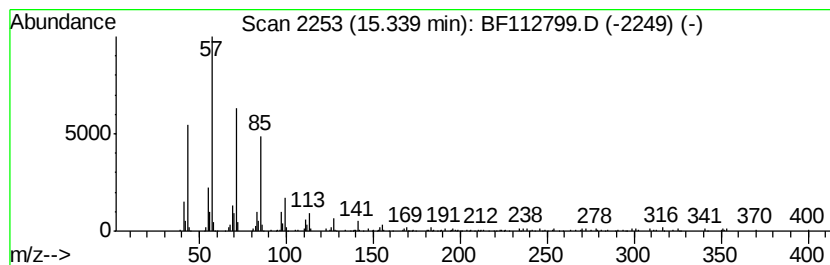
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 14 Octadecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.34	3.82 ng	207037	Perylene-d12	15.28

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecane	254	C18H38	000593-45-3	91
2		Heneicosane	296	C21H44	000629-94-7	90
3		Nonadecane	268	C19H40	000629-92-5	86
4		Hentriacontane	437	C31H64	000630-04-6	86
5		Docosane	310	C22H46	000629-97-0	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF030119\
Data File : BF112799.D
Acq On : 1 Mar 2019 14:58
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Sample : K1655-02
Misc :
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Instrument :
BNA_F
ClientSampleId :
DUP-2

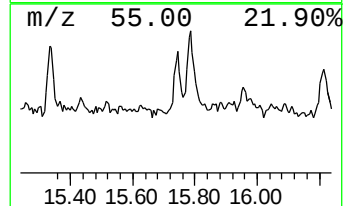
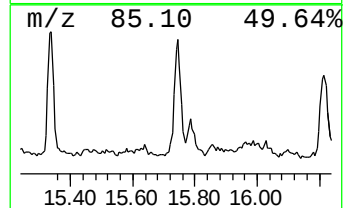
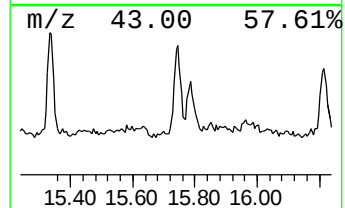
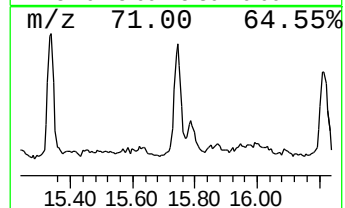
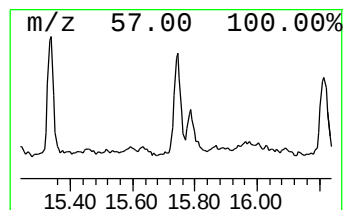
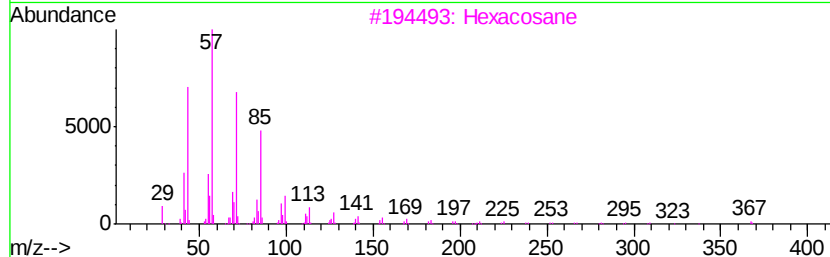
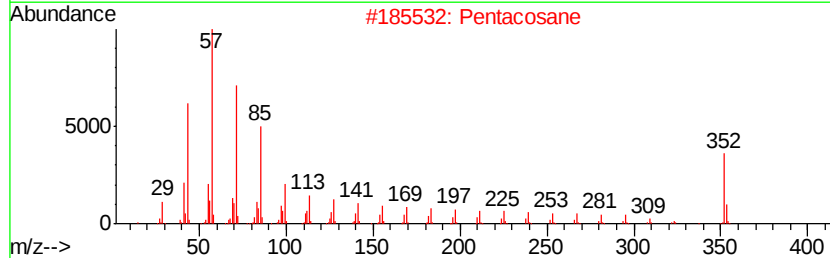
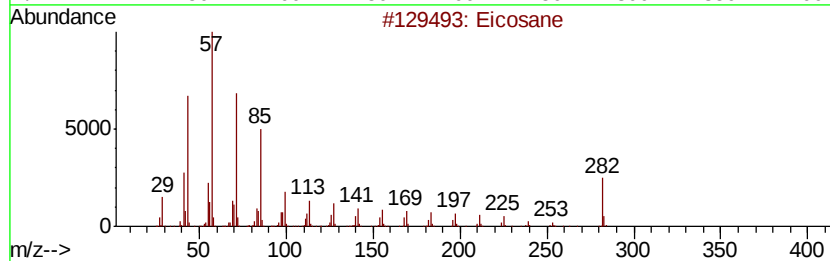
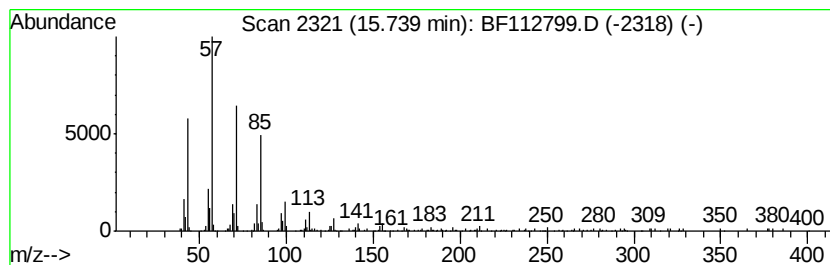
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Eicosane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.74	4.17 ng	225893	Perylene-d12	15.28

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	91
2		Pentacosane	352	C25H52	000629-99-2	91
3		Hexacosane	366	C26H54	000630-01-3	87
4		Eicosane, 7-hexyl-	366	C26H54	055333-99-8	86
5		Nonadecane, 2-methyl-	282	C20H42	001560-86-7	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF030119\
Data File : BF112799.D
Acq On : 1 Mar 2019 14:58
Operator : JU/SJ
Sample : K1655-02
Misc :
ALS Vial : 10 Sample Multiplier: 1

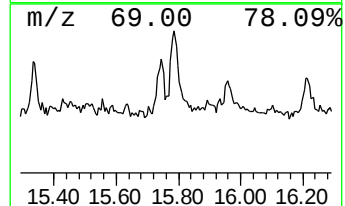
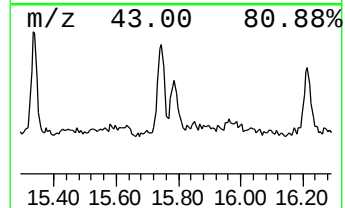
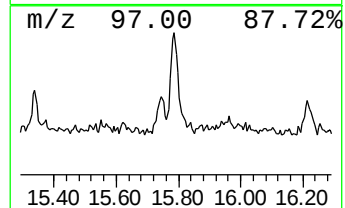
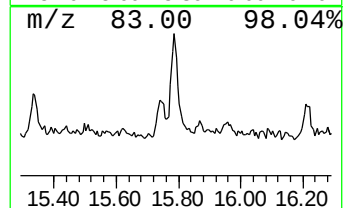
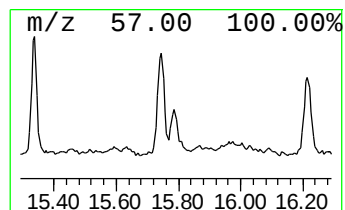
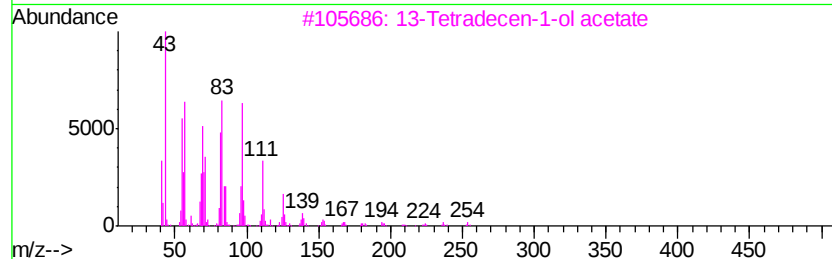
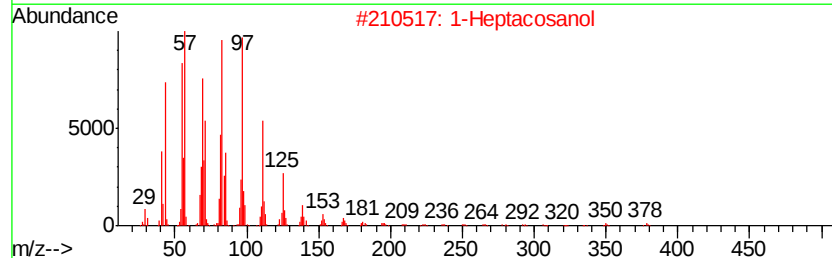
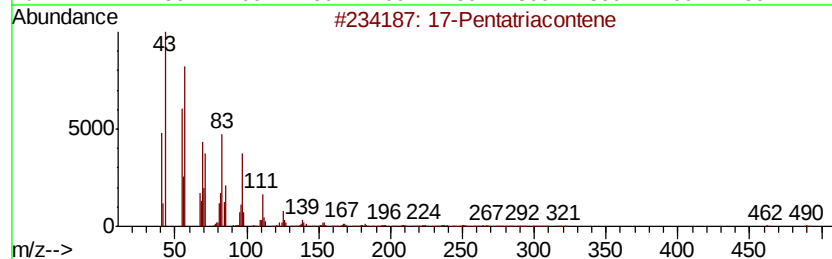
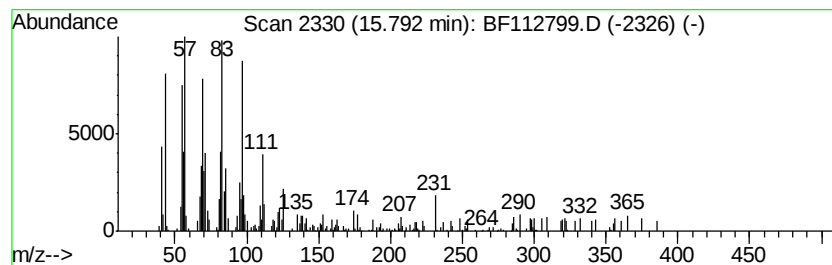
Instrument :
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 17-Pentatriacontene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.79	4.10 ng	222154	Perylene-d12	15.28	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	17-Pentatriacontene	491	C35H70	006971-40-0	58
2	1-Heptacosanol	396	C27H56O	002004-39-9	58
3	13-Tetradecen-1-ol acetate	254	C16H30O2	056221-91-1	53
4	Octacosanol	410	C28H58O	000557-61-9	53
5	Cyclopentane, 1,1,3-trimethyl-	112	C8H16	004516-69-2	49



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF030119\
Data File : BF112799.D
Acq On : 1 Mar 2019 14:58
Operator : JU/SJ
Sample : K1655-02
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
DUP-2

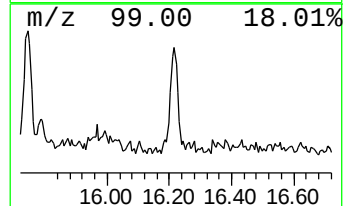
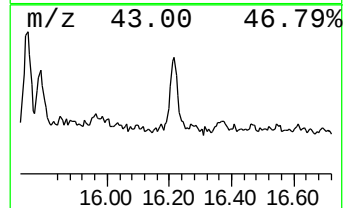
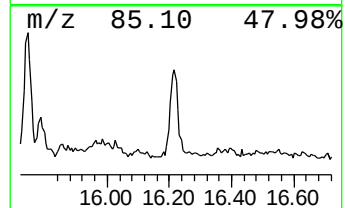
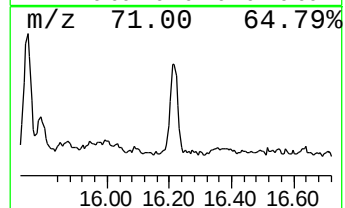
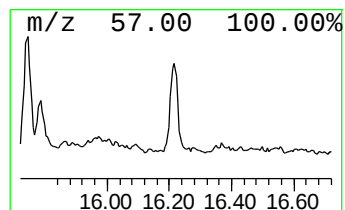
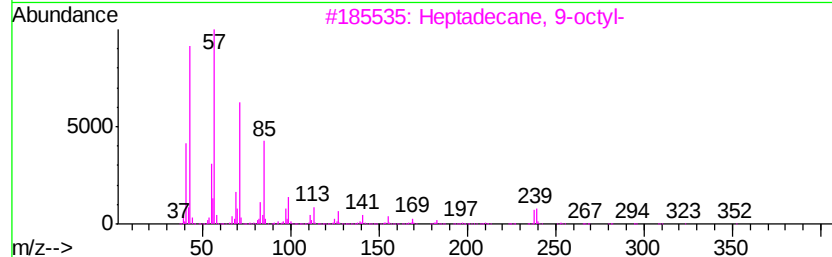
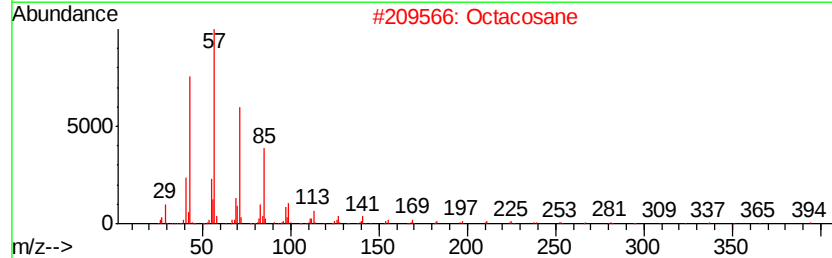
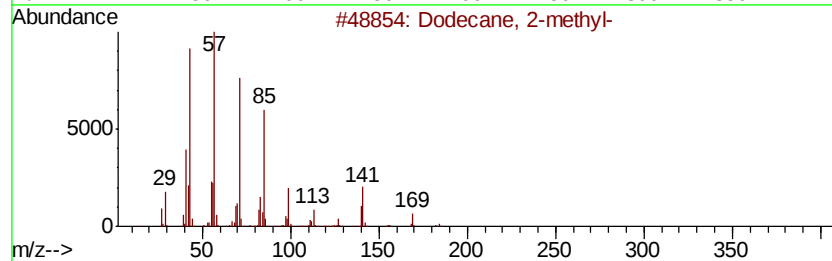
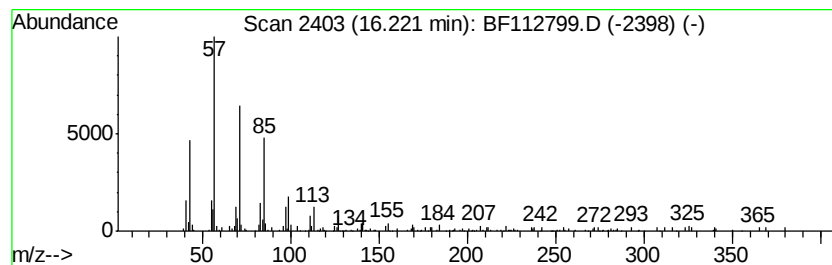
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 Dodecane, 2-methyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.22	3.09 ng	167701	Perylene-d12	15.28

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecane, 2-methyl-	184	C13H28	001560-97-0	90
2			Octacosane	394	C28H58	000630-02-4	87
3			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	87
4			Nonadecane	268	C19H40	000629-92-5	83
5			Hexadecane	226	C16H34	000544-76-3	80



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF030119\
 Data File : BF112799.D
 Acq On : 1 Mar 2019 14:58
 Operator : JU/SJ
 Sample : K1655-02
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF022719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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Hexadecanoic acid...	11.66	3.3	ng	160031	4	11.24	961616	20.0
n-Hexadecanoic acid	11.79	14.3	ng	689701	4	11.24	961616	20.0
Methyl stearate	12.44	2.5	ng	122581	4	11.24	961616	20.0
Octadecanoic acid	12.57	6.3	ng	352140	5	13.87	1123830	20.0
Hexadecanoic acid...	12.66	3.0	ng	167041	5	13.87	1123830	20.0
Octadecanoic acid...	13.37	3.6	ng	200205	5	13.87	1123830	20.0
1-Heneicosanol	13.73	9.9	ng	556599	5	13.87	1123830	20.0
Heptadecane	14.06	2.4	ng	136801	5	13.87	1123830	20.0
Heptacosane, 1-ch...	14.36	3.4	ng	192179	5	13.87	1123830	20.0
Heneicosane	14.66	4.0	ng	216027	6	15.28	1084030	20.0
1-Iodo-2-methylun...	14.98	3.9	ng	209443	6	15.28	1084030	20.0
Octadecane	15.34	3.8	ng	207037	6	15.28	1084030	20.0
Eicosane	15.74	4.2	ng	225893	6	15.28	1084030	20.0
17-Pentatriacontene	15.79	4.1	ng	222154	6	15.28	1084030	20.0
Dodecane, 2-methyl-	16.22	3.1	ng	167701	6	15.28	1084030	20.0