

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF021919\
 Data File : BF112622.D
 Acq On : 19 Feb 2019 15:46
 Operator : JU/SJ
 Sample : K1515-09
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TS-9

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-BF012519.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	5.051	500	504	521	rBV	75581	112727	3.83%	0.444%
2	5.463	570	574	603	rBV	2000430	2341420	79.56%	9.224%
3	6.475	742	746	763	rBV	2299592	2624117	89.16%	10.338%
4	6.598	763	767	781	rVB	2792780	2676184	90.93%	10.543%
5	6.822	801	805	814	rBV	566034	511301	17.37%	2.014%
6	6.981	827	832	844	rBV	1974401	1992839	67.71%	7.851%
7	7.386	897	901	922	rBV	1527505	1602615	54.45%	6.314%
8	8.104	1019	1023	1041	rBV	656394	662700	22.52%	2.611%
9	9.175	1200	1205	1216	rBV	2984691	2943125	100.00%	11.595%
10	9.575	1269	1273	1279	rBV	133012	141408	4.80%	0.557%
11	9.857	1317	1321	1330	rVB	824573	786835	26.73%	3.100%
12	10.657	1453	1457	1471	rBV	2070268	2096305	71.23%	8.259%
13	11.351	1571	1575	1583	rBV	749470	816683	27.75%	3.217%
14	11.868	1659	1663	1674	rBV2	241351	286451	9.73%	1.129%
15	12.574	1778	1783	1788	rBV	20589	35881	1.22%	0.141%
16	12.651	1793	1796	1807	rBV	68724	100387	3.41%	0.395%
17	12.927	1839	1843	1846	rBV	3008200	2802651	95.23%	11.041%
18	13.145	1874	1880	1883	rBV3	28974	36565	1.24%	0.144%
19	13.486	1934	1938	1941	rBV2	32122	33528	1.14%	0.132%
20	13.815	1990	1994	2005	rBV	196691	288479	9.80%	1.137%
21	13.998	2021	2025	2032	rBV	653597	693141	23.55%	2.731%
22	14.127	2044	2047	2050	rBV	91556	74832	2.54%	0.295%
23	14.433	2095	2099	2101	rBV	145066	141187	4.80%	0.556%
24	14.733	2146	2150	2155	rVB	165464	151782	5.16%	0.598%
25	14.804	2159	2162	2166	rBV	131063	133891	4.55%	0.527%
26	15.062	2200	2206	2211	rBV	178826	215104	7.31%	0.847%
27	15.433	2264	2269	2272	rBV	127431	175369	5.96%	0.691%
28	15.468	2272	2275	2285	rVB	440869	584296	19.85%	2.302%
29	15.862	2336	2342	2347	rBV	107249	159790	5.43%	0.630%
30	16.351	2420	2425	2430	rBV	65145	104209	3.54%	0.411%
31	16.933	2520	2524	2530	rVB4	31109	57271	1.95%	0.226%

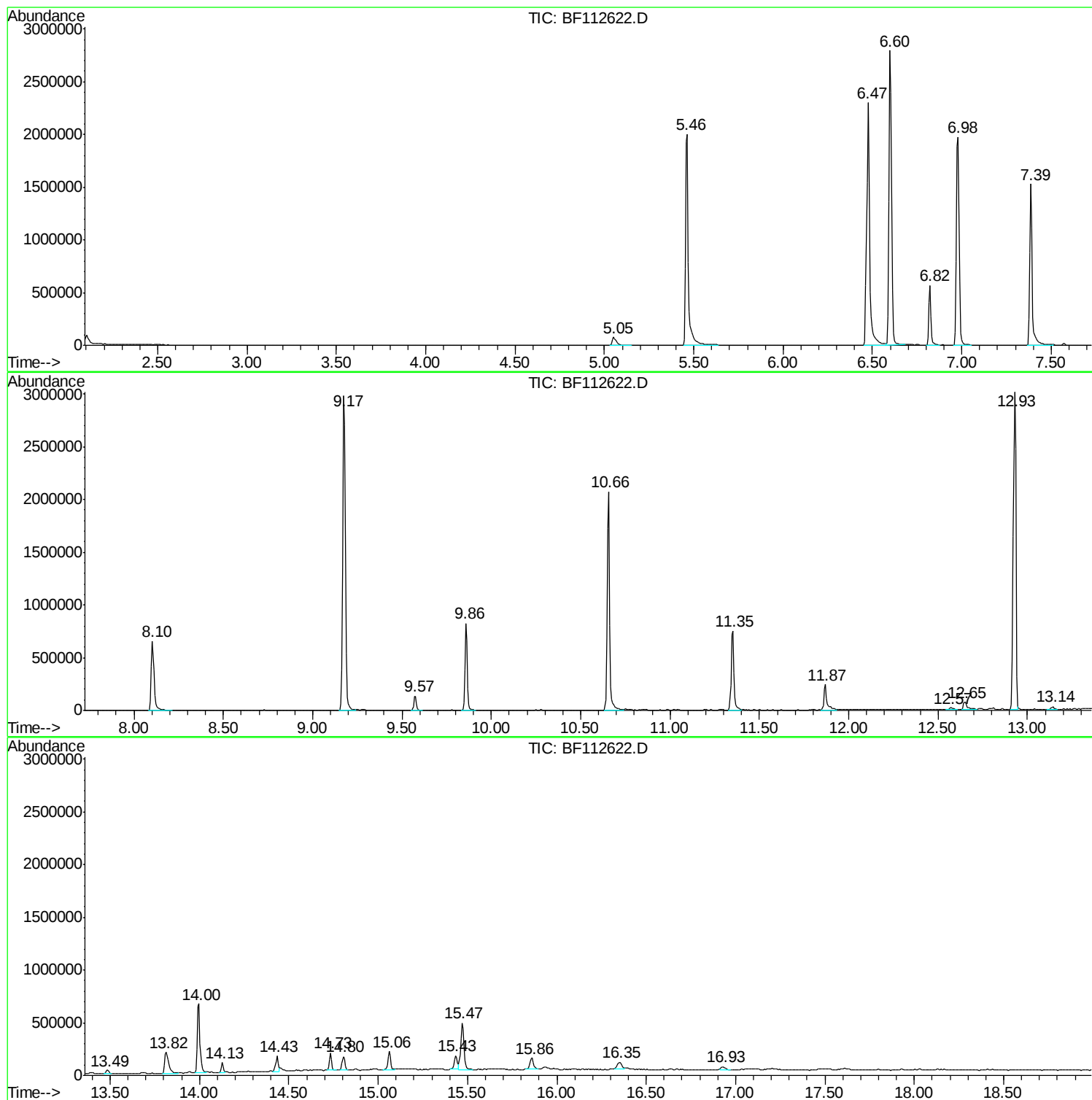
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF012519.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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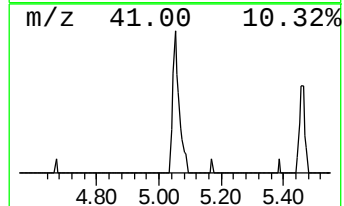
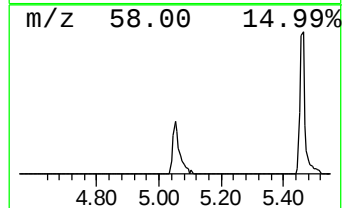
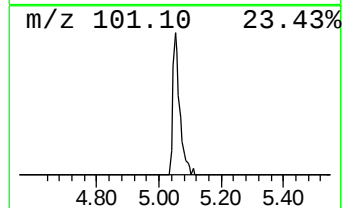
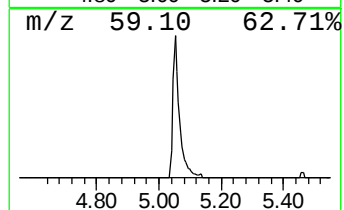
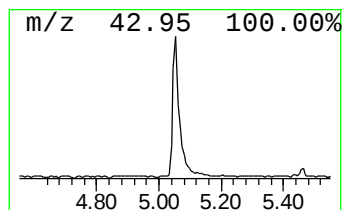
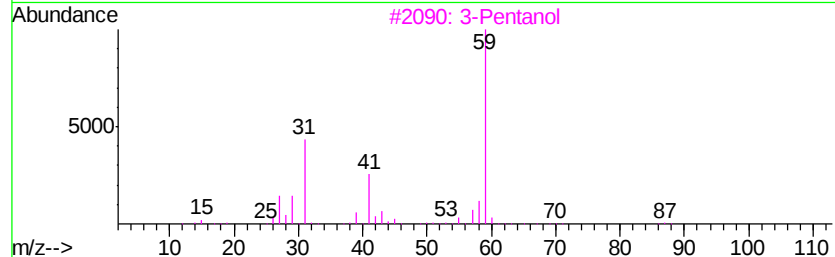
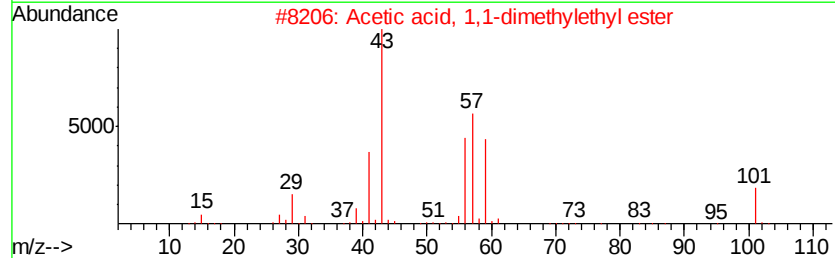
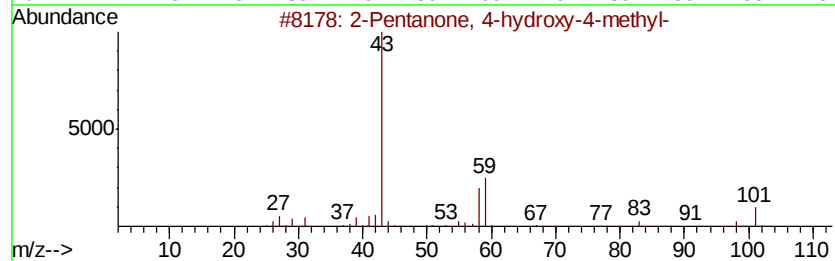
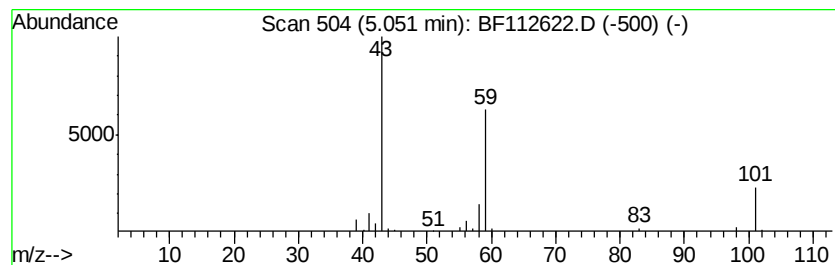
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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.
5.05	4.41 ng	112727	1,4-Dichlorobenzene-d4	6.82

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3			3-Pentanol	88	C5H12O	000584-02-1	37
4			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25
5			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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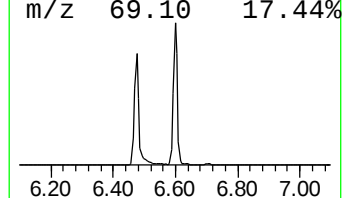
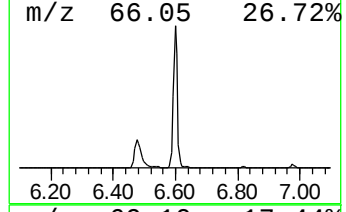
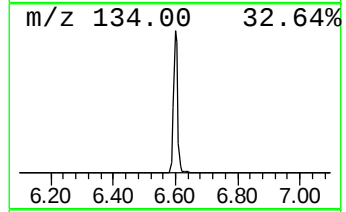
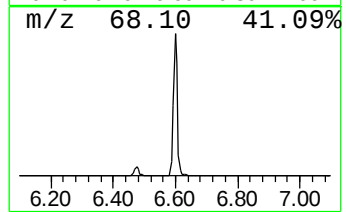
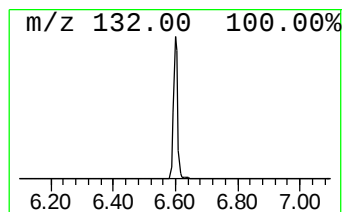
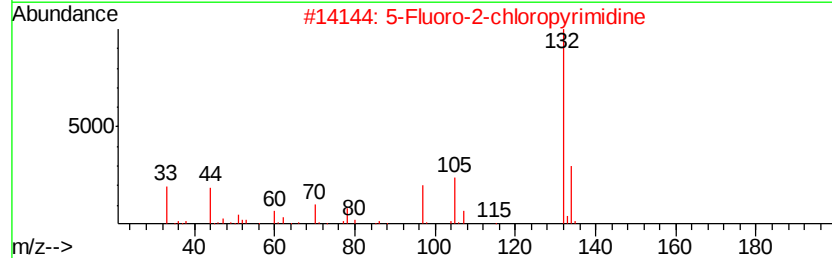
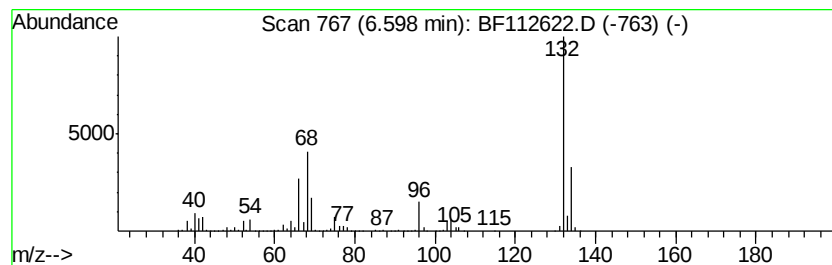
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown6.60 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.60	104.68 ng	2676180	1,4-Dichlorobenzene-d4	6.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
3		Tranvlcyproline-propionyl	189	C12H15NO	1000123-86-3	10
4		1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9
5		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9



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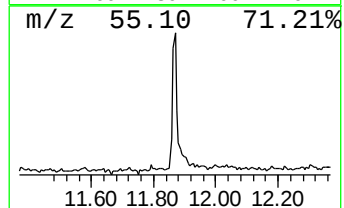
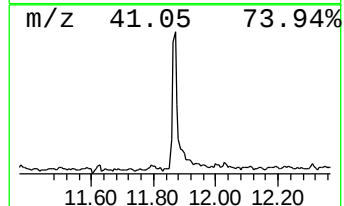
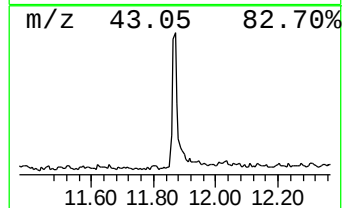
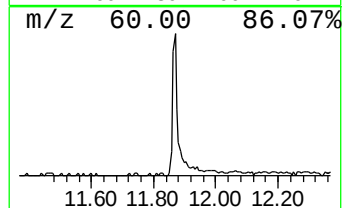
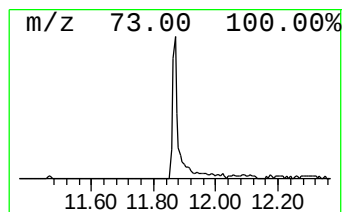
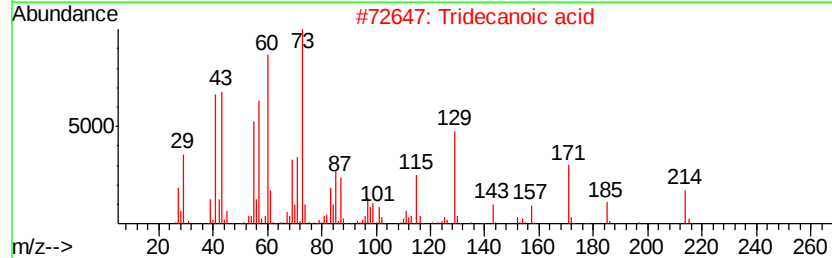
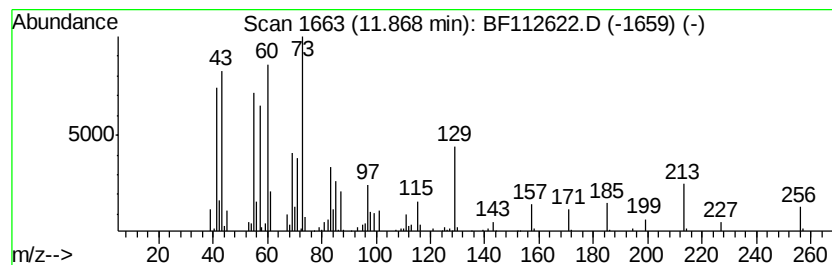
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 Peak Number 4 n-Hexadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD		R.T.	
11.87	7.01 ng	286451	Phenanthrene-d10		11.35	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		Tridecanoic acid	214	C13H26O2	000638-53-9	91
3		Pentadecanoic acid	242	C15H30O2	001002-84-2	90
4		Undecanoic acid	186	C11H22O2	000112-37-8	80
5		Tetradecanoic acid	228	C14H28O2	000544-63-8	80



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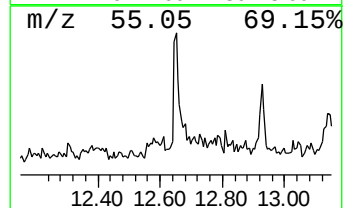
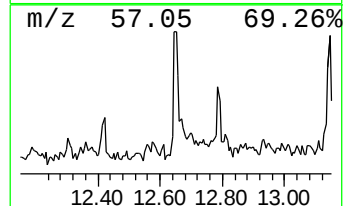
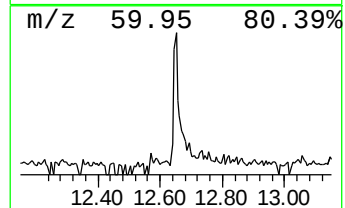
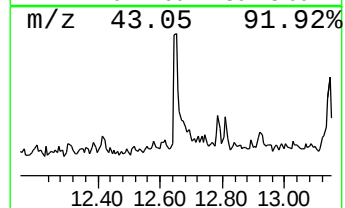
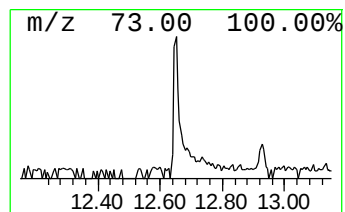
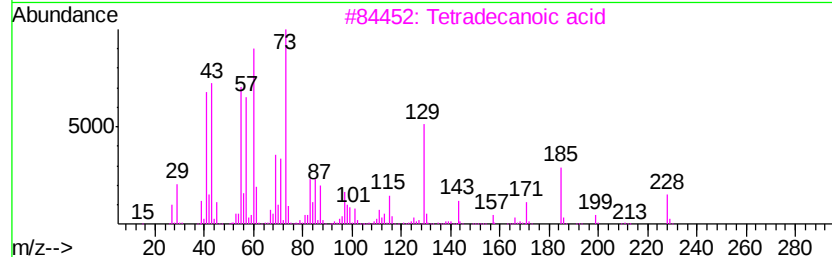
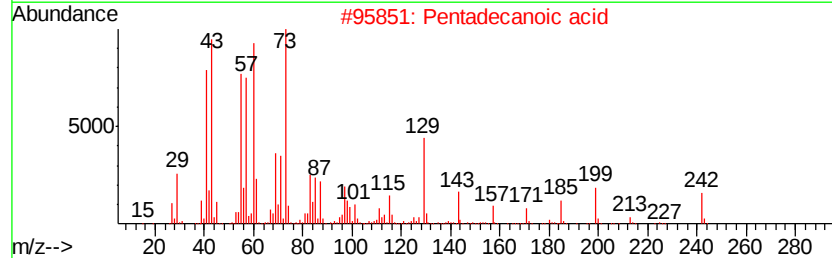
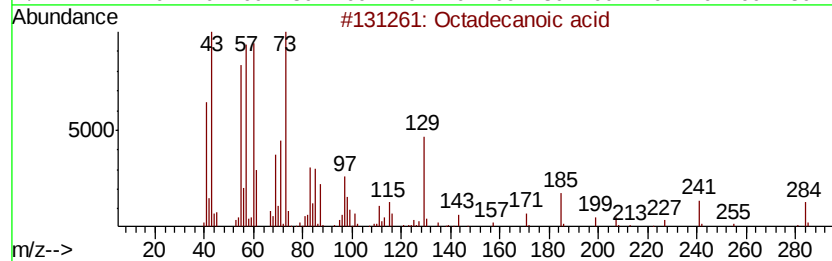
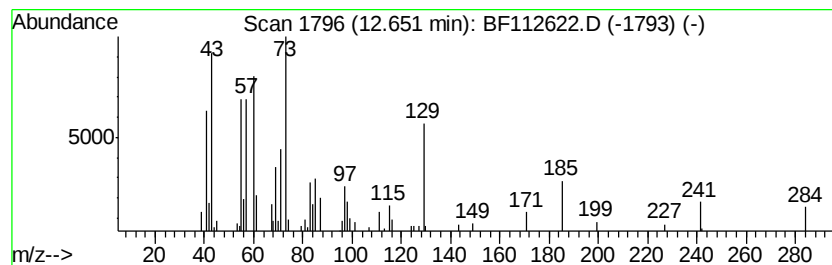
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 5 Octadecanoic acid Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD		R.T.
12.65	2.46 ng	100387	Phenanthrene-d10		11.35
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	Tetradecanoic acid	228	C14H28O2	000544-63-8	87
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	72
5	n-Decanoic acid	172	C10H20O2	000334-48-5	59



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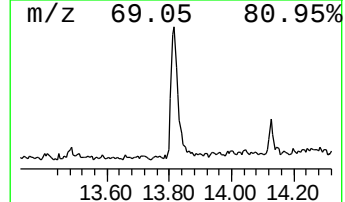
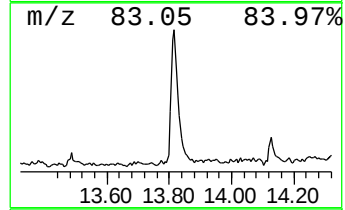
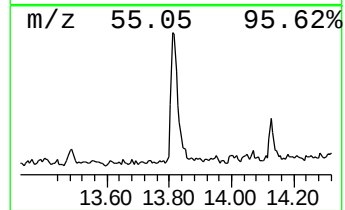
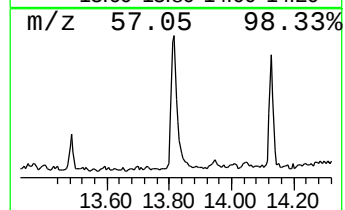
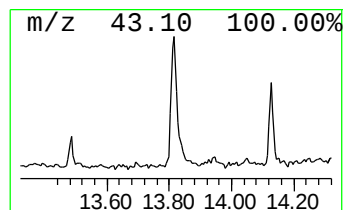
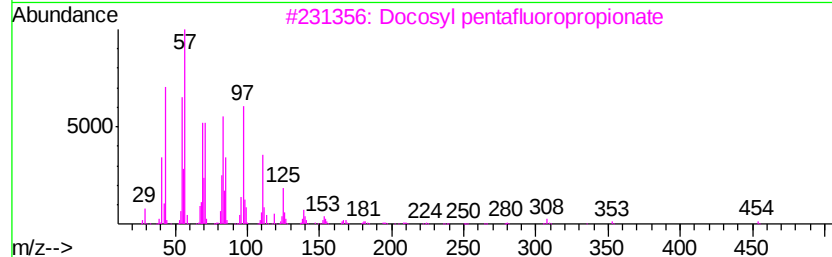
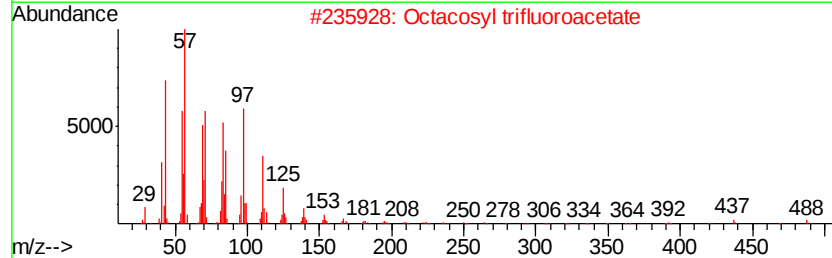
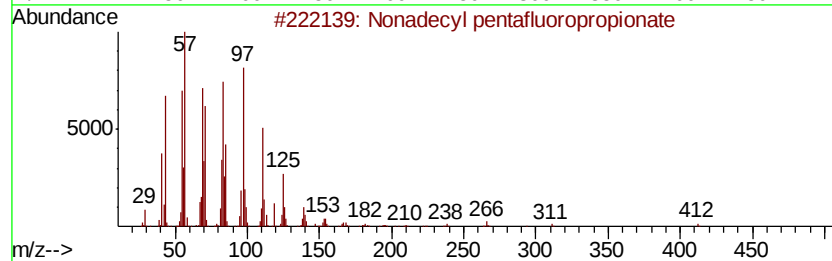
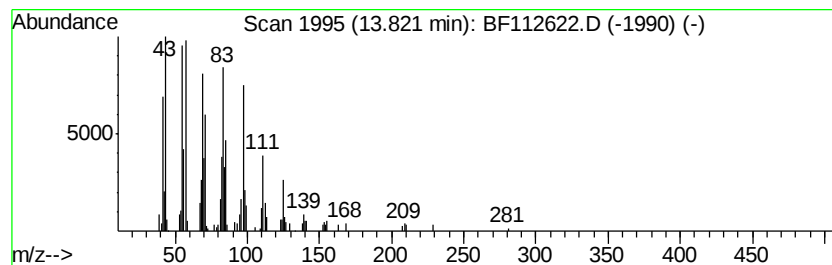
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 6 Nonadecyl pentafluoropropio... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.82	8.32 ng	288479	Chrysene-d12	14.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonadecyl pentafluoropropionate	430	C22H39F5O2	1000351-88-8	91
2		Octacosyl trifluoroacetate	506	C30H57F3O2	1000351-74-9	91
3		Docosyl pentafluoropropionate	472	C25H45F5O2	1000351-80-9	91
4		1-Dodecanol, 2-hexyl-	270	C18H38O	110225-00-8	90
5		17-Pentatriacontene	491	C35H70	006971-40-0	83



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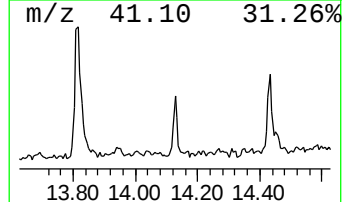
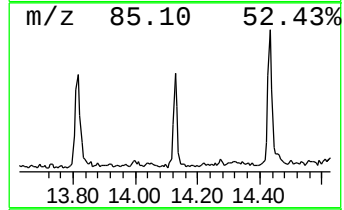
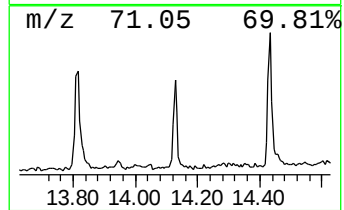
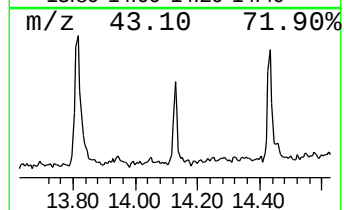
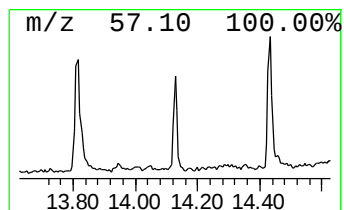
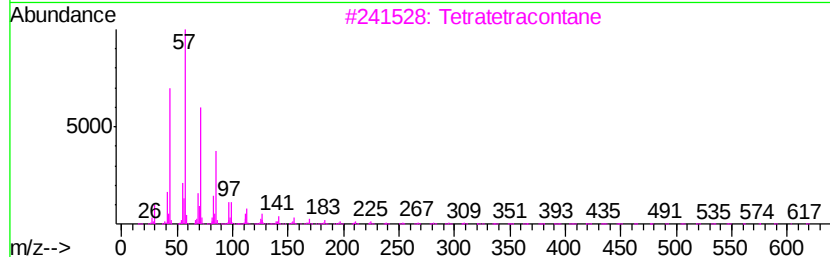
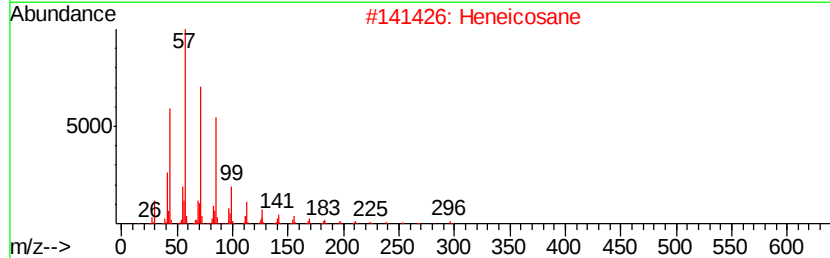
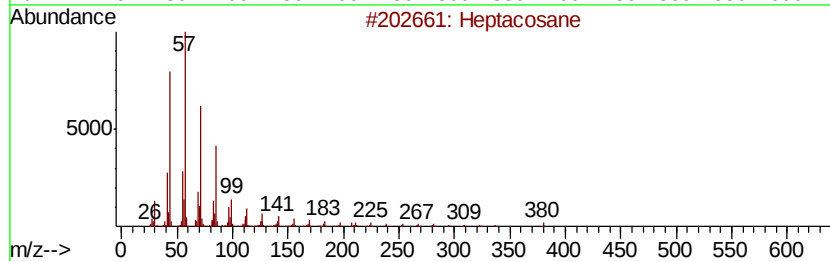
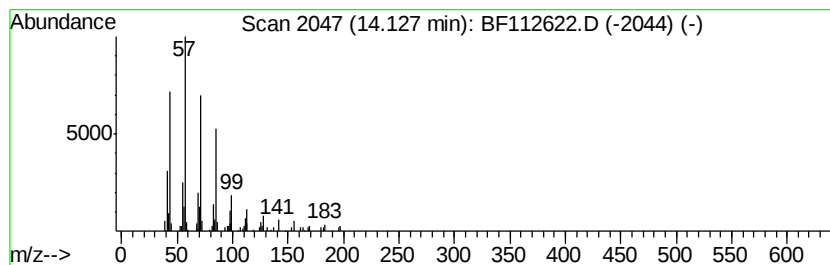
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Heptacosane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.13	2.16 ng	74832	Chrysene-d12	14.00

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptacosane		380	C27H56	000593-49-7	91
2	Heneicosane		296	C21H44	000629-94-7	91
3	Tetratetracontane		619	C44H90	007098-22-8	91
4	Hexacosane		366	C26H54	000630-01-3	91
5	Tetracosane		338	C24H50	000646-31-1	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF021919\
Data File : BF112622.D
Acq On : 19 Feb 2019 15:46
Operator : JU/SJ
Sample : K1515-09
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TS-9

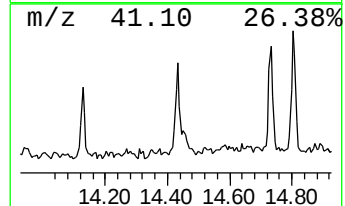
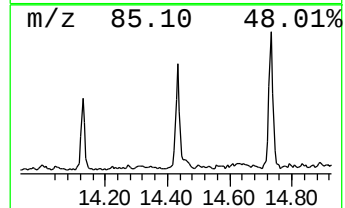
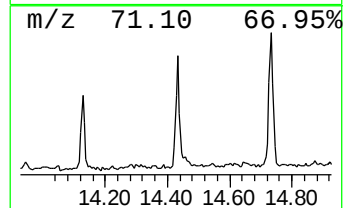
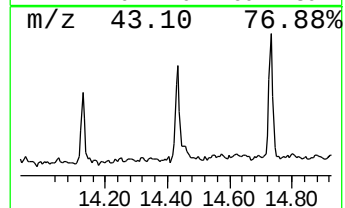
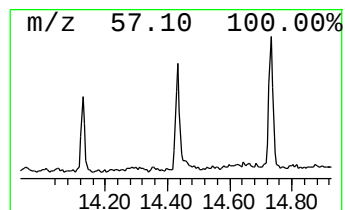
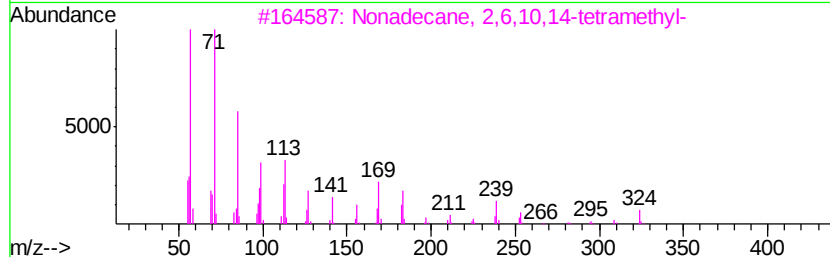
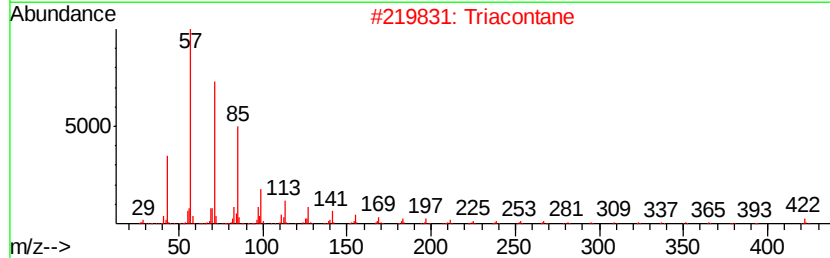
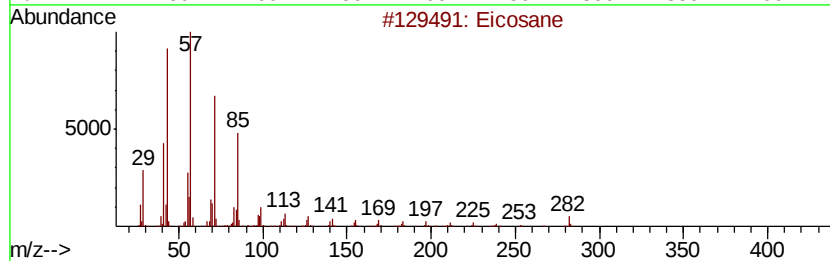
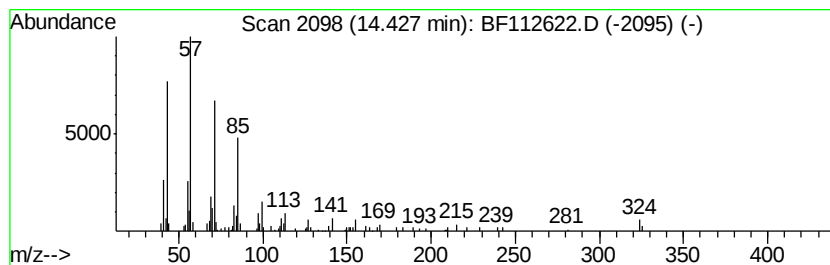
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Eicosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.43	4.07 ng	141187	Chrysene-d12	14.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	96
2		Triacontane	422	C30H62	000638-68-6	90
3		Nonadecane, 2,6,10,14-tetramethyl-	324	C23H48	055124-80-6	90
4		Heneicosane	296	C21H44	000629-94-7	90
5		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF021919\
Data File : BF112622.D
Acq On : 19 Feb 2019 15:46
Operator : JU/SJ
Sample : K1515-09
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TS-9

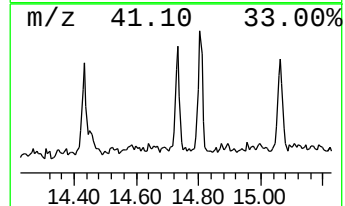
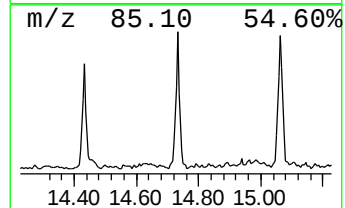
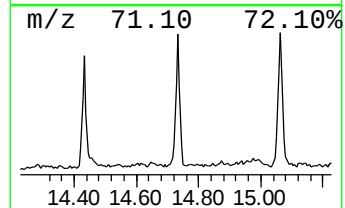
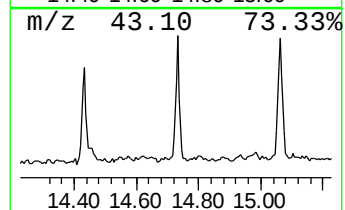
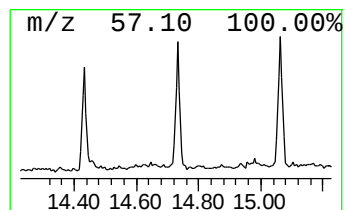
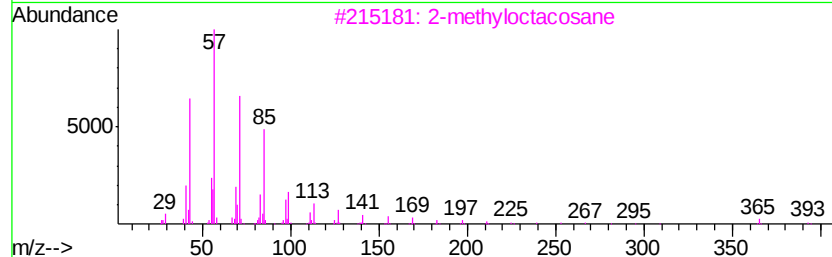
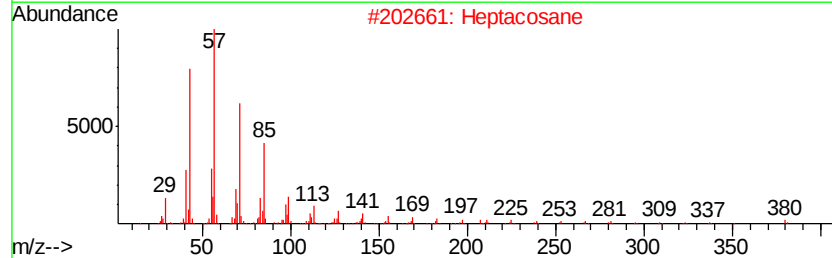
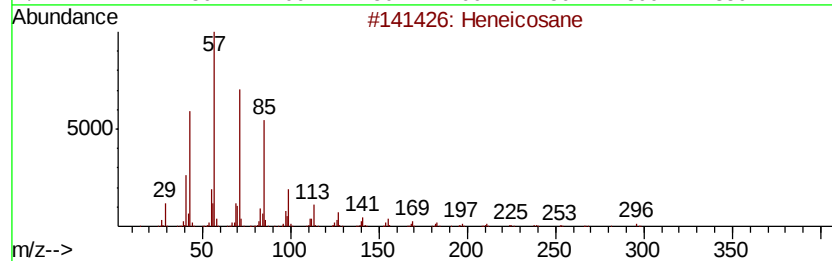
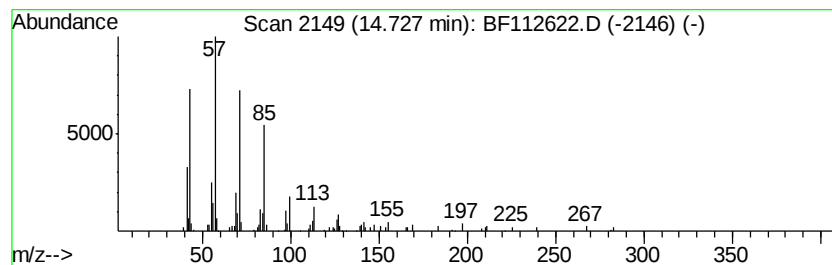
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Heneicosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.73	4.38 ng	151782	Chrysene-d12	14.00

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane		296	C21H44	000629-94-7	91
2	Heptacosane		380	C27H56	000593-49-7	90
3	2-methyloctacosane		408	C29H60	1000376-72-8	90
4	Hexacosane		366	C26H54	000630-01-3	90
5	5-Ethyl-5-methylnonadecane		310	C22H46	1000360-42-7	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF021919\
Data File : BF112622.D
Acq On : 19 Feb 2019 15:46
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Sample : K1515-09
Misc :
ALS Vial : 12 Sample Multiplier: 1

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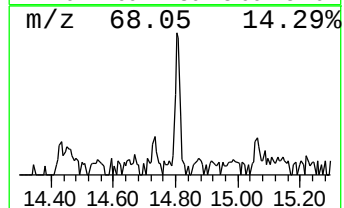
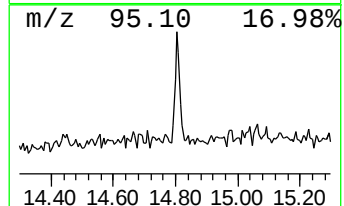
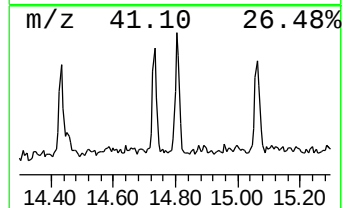
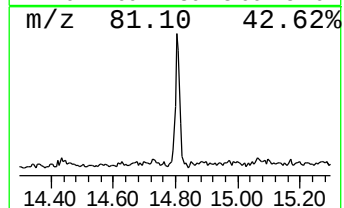
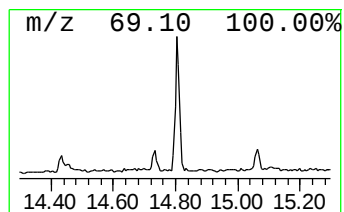
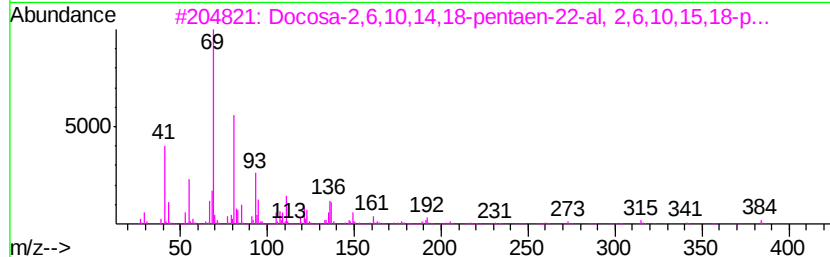
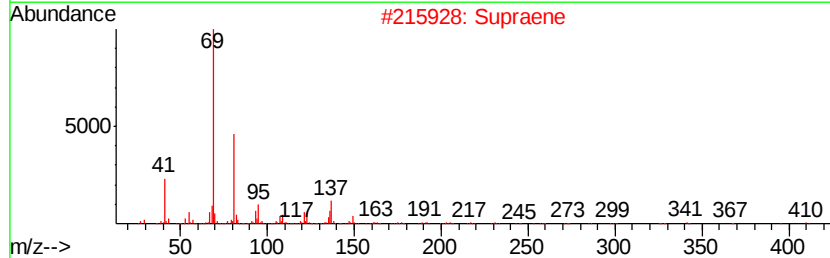
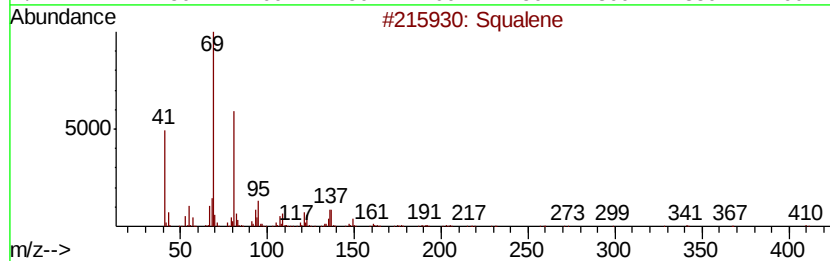
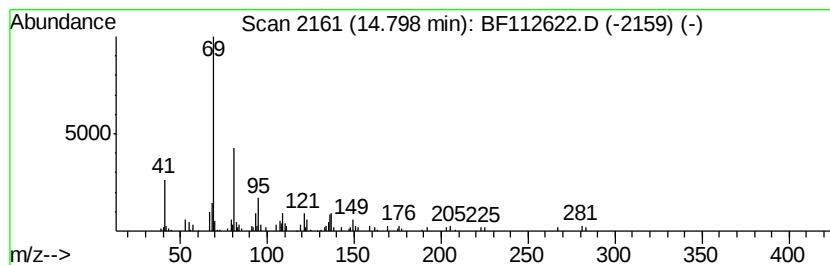
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Squalene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.80	4.58 ng	133891	Perylene-d12	15.47

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Squalene	410	C30H50	000111-02-4	91
2		Supraene	410	C30H50	007683-64-9	91
3		Docosa-2,6,10,14,18-pentaen-22-a...	384	C27H44O	1000163-04-7	83
4		7,11-Dimethyldodeca-2,6,10-trien...	208	C14H24O	125529-10-4	74
5		Propanoic acid, 2,2-dimethyl-, [...]	306	C20H34O2	1000164-38-8	59



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF021919\
Data File : BF112622.D
Acq On : 19 Feb 2019 15:46
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Sample : K1515-09
Misc :
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ClientSampleId :
TS-9

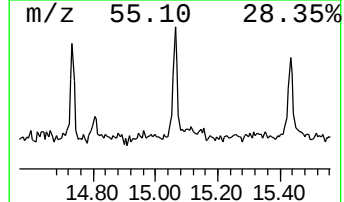
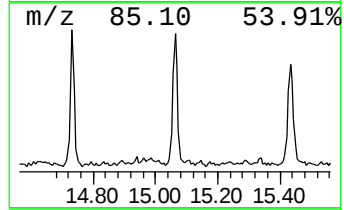
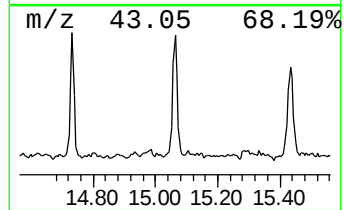
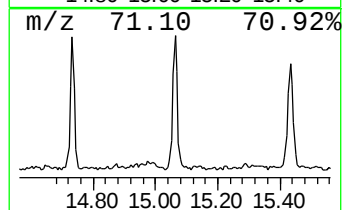
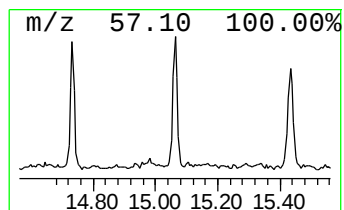
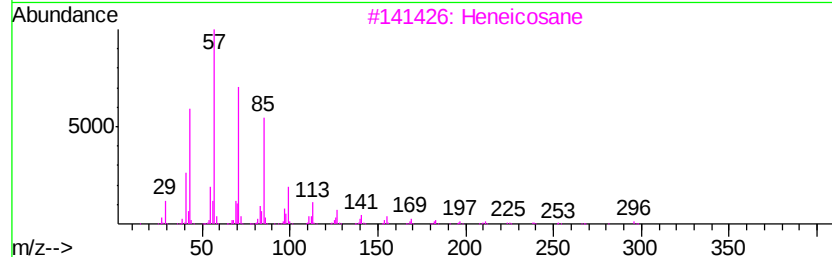
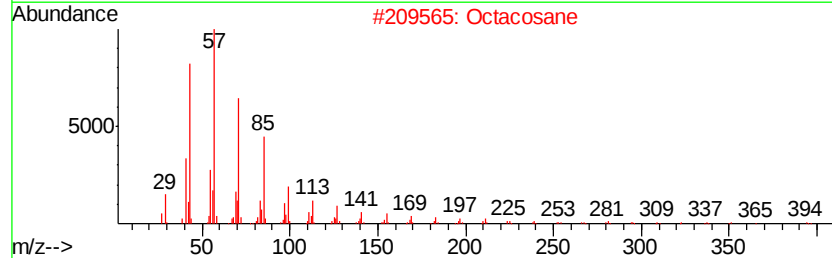
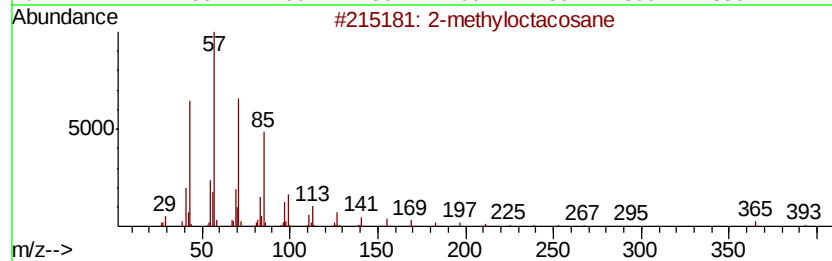
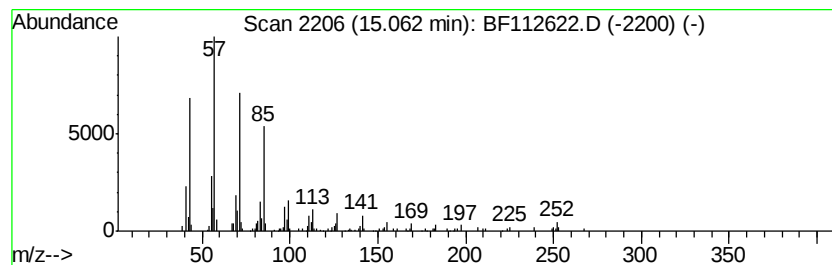
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 2-methyloctacosane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.06	7.36 ng	215104	Perylene-d12	15.47

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-methyloctacosane	408	C29H60	1000376-72-8	90
2		Octacosane	394	C28H58	000630-02-4	90
3		Heneicosane	296	C21H44	000629-94-7	90
4		Hexacosane	366	C26H54	000630-01-3	90
5		Octadecane	254	C18H38	000593-45-3	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF021919\
Data File : BF112622.D
Acq On : 19 Feb 2019 15:46
Operator : JU/SJ
Sample : K1515-09
Misc :
ALS Vial : 12 Sample Multiplier: 1

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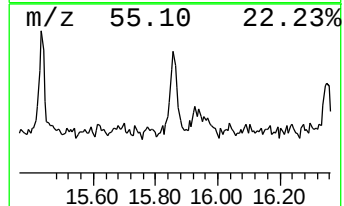
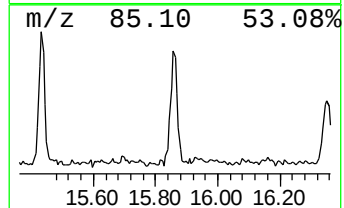
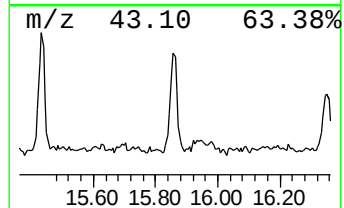
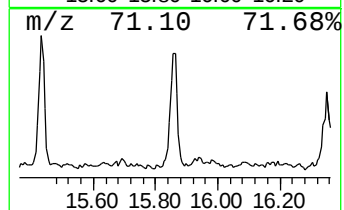
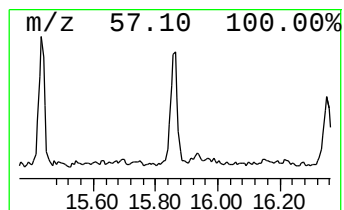
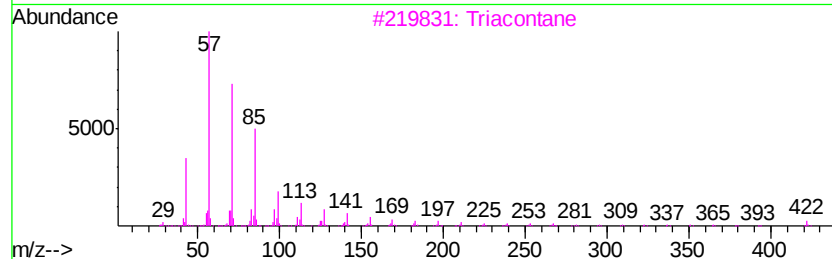
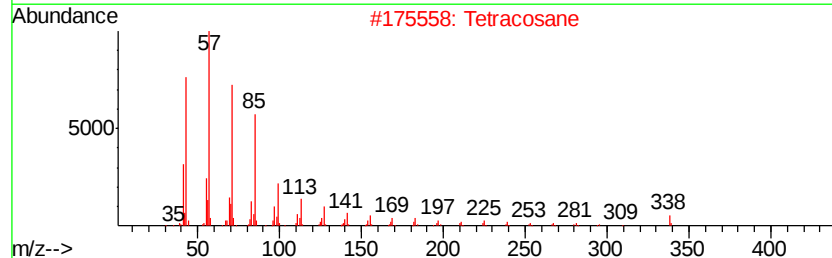
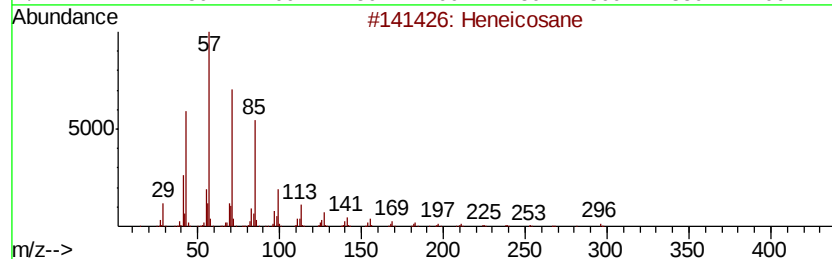
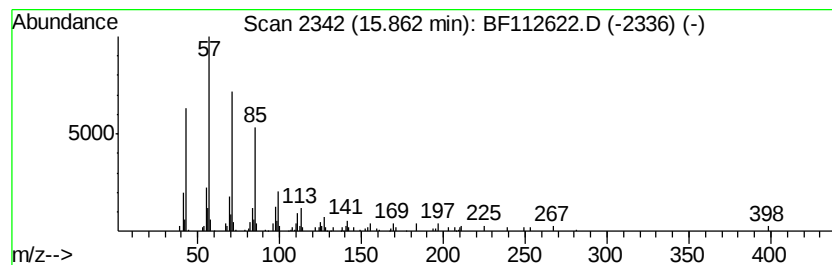
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Sulfurous acid, butyl dodec... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.86	5.47 ng	159790	Perylene-d12	15.47

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	87
2		Tetracosane	338	C24H50	000646-31-1	87
3		Triacotane	422	C30H62	000638-68-6	87
4		Hexacosane	366	C26H54	000630-01-3	83
5		Sulfurous acid, butyl dodecyl ester	306	C16H34O3S	1000309-17-9	80



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF021919\
Data File : BF112622.D
Acq On : 19 Feb 2019 15:46
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Sample : K1515-09
Misc :
ALS Vial : 12 Sample Multiplier: 1

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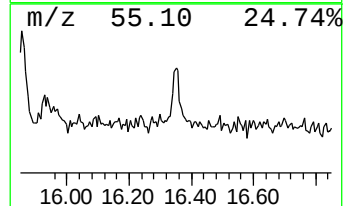
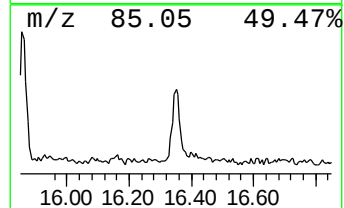
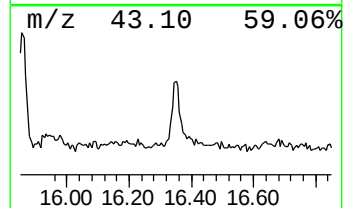
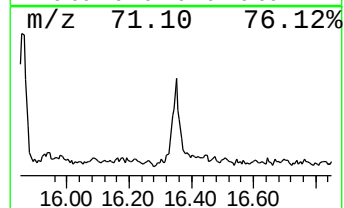
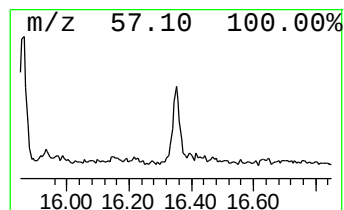
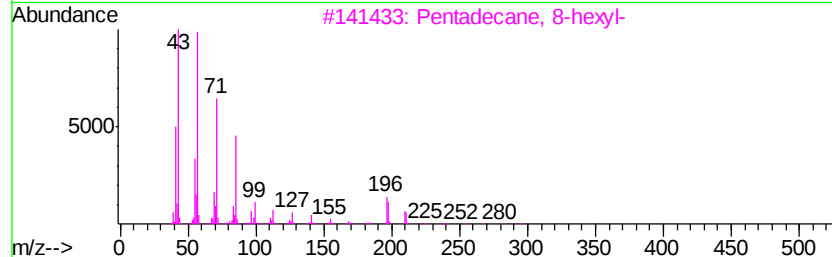
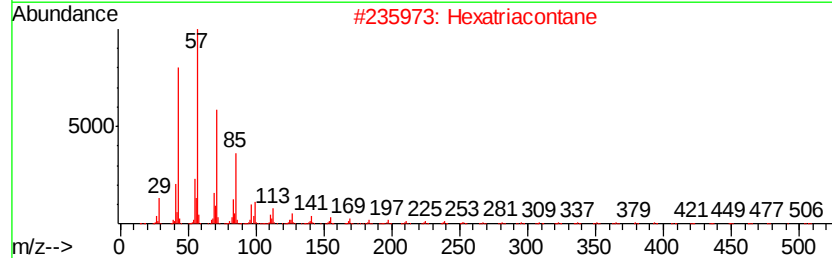
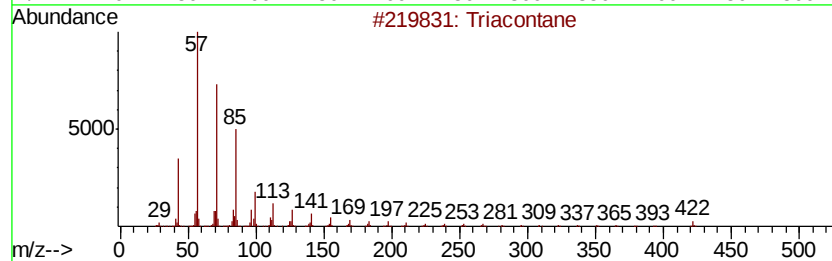
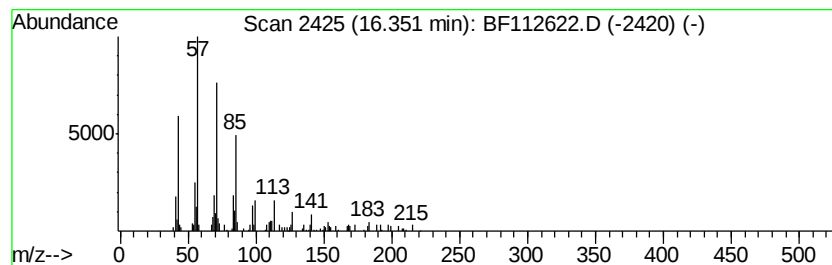
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Triacontane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.35	3.57 ng	104209	Perylene-d12	15.47

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Triacontane	422	C30H62	000638-68-6	90
2			Hexatriacontane	507	C36H74	000630-06-8	90
3			Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	90
4			Pentacosane	352	C25H52	000629-99-2	90
5			Hexacosane	366	C26H54	000630-01-3	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF021919\
 Data File : BF112622.D
 Acq On : 19 Feb 2019 15:46
 Operator : JU/SJ
 Sample : K1515-09
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF012519.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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unknown6.60	6.60	104.7	ng	2676180	1	6.82	511301	20.0
n-Hexadecanoic acid	11.87	7.0	ng	286451	4	11.35	816683	20.0
Octadecanoic acid	12.65	2.5	ng	100387	4	11.35	816683	20.0
Nonadecyl pentafl...	13.82	8.3	ng	288479	5	14.00	693141	20.0
Heptacosane	14.13	2.2	ng	74832	5	14.00	693141	20.0
Eicosane	14.43	4.1	ng	141187	5	14.00	693141	20.0
Heneicosane	14.73	4.4	ng	151782	5	14.00	693141	20.0
Squalene	14.80	4.6	ng	133891	6	15.47	584296	20.0
2-methyloctacosane	15.06	7.4	ng	215104	6	15.47	584296	20.0
Sulfurous acid, b...	15.86	5.5	ng	159790	6	15.47	584296	20.0
Triacontane	16.35	3.6	ng	104209	6	15.47	584296	20.0