

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF020619\
 Data File : BF112419.D
 Acq On : 6 Feb 2019 14:52
 Operator : JU/SJ
 Sample : K1338-03 5X
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 AU-06-020519-B

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 3 % of largest Peak

Max Peaks: 100

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-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	4.557	415	420	425	rVB4	10004	14364	1.08%	0.099%
2	5.040	499	502	506	rBV	17661	20259	1.53%	0.139%
3	5.463	571	574	590	rBV	406504	505944	38.09%	3.475%
4	6.481	743	747	765	rBV	408316	558458	42.04%	3.836%
5	6.604	765	768	778	rBV	603826	600717	45.22%	4.126%
6	6.828	803	806	816	rBV	860199	828326	62.36%	5.690%
7	6.987	829	833	843	rBV	519604	462687	34.83%	3.178%
8	7.392	899	902	913	rBV	314402	344003	25.90%	2.363%
9	8.110	1020	1024	1033	rBV	1078297	1097434	82.62%	7.538%
10	8.169	1033	1034	1038	rVB	19135	16458	1.24%	0.113%
11	8.486	1085	1088	1093	rBV4	15505	16376	1.23%	0.112%
12	8.534	1093	1096	1098	rBV	34585	30011	2.26%	0.206%
13	9.186	1201	1207	1218	rBV	728734	668465	50.32%	4.592%
14	9.581	1271	1274	1282	rBV	37150	43414	3.27%	0.298%
15	9.798	1303	1311	1313	rBV4	11087	18428	1.39%	0.127%
16	9.869	1318	1323	1334	rVB	1452447	1310993	98.70%	9.005%
17	10.657	1454	1457	1470	rBV	401010	439859	33.11%	3.021%
18	10.769	1473	1476	1483	rVB4	9815	18494	1.39%	0.127%
19	11.357	1572	1576	1588	rBV	1479891	1328317	100.00%	9.124%
20	11.739	1637	1641	1645	rBV	400628	322834	24.30%	2.218%
21	11.874	1659	1664	1672	rBV4	38982	66301	4.99%	0.455%
22	12.422	1753	1757	1759	rBV	1151109	1020153	76.80%	7.007%
23	12.445	1759	1761	1764	rVB	1305869	994683	74.88%	6.832%
24	12.527	1772	1775	1780	rVB	213527	173710	13.08%	1.193%
25	12.574	1780	1783	1787	rBV	33475	39294	2.96%	0.270%
26	12.663	1795	1798	1803	rBV5	12082	17168	1.29%	0.118%
27	12.804	1819	1822	1824	rBV	35736	32133	2.42%	0.221%
28	12.933	1840	1844	1852	rVB	690254	610213	45.94%	4.192%
29	13.257	1896	1899	1903	rVB4	13658	14212	1.07%	0.098%
30	13.504	1938	1941	1943	rVB3	18959	16866	1.27%	0.116%
31	13.827	1992	1996	1999	rBV	68938	80817	6.08%	0.555%
32	13.921	2010	2012	2016	rBV4	14896	16786	1.26%	0.115%
33	13.998	2021	2025	2035	rBV	1014019	966910	72.79%	6.642%
34	14.145	2047	2050	2053	rBV	57747	46032	3.47%	0.316%

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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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Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	14.451	2099	2102	2106	rBV	77277	70357	5.30%	0.483%
36	14.751	2150	2153	2157	rVB	95849	87782	6.61%	0.603%
37	15.080	2206	2209	2216	rVB	134985	181780	13.68%	1.249%
38	15.468	2269	2275	2282	rBV	802560	1092539	82.25%	7.505%
39	15.886	2342	2346	2352	rBV	147848	217379	16.36%	1.493%
40	16.386	2426	2431	2437	rVB2	93132	167261	12.59%	1.149%

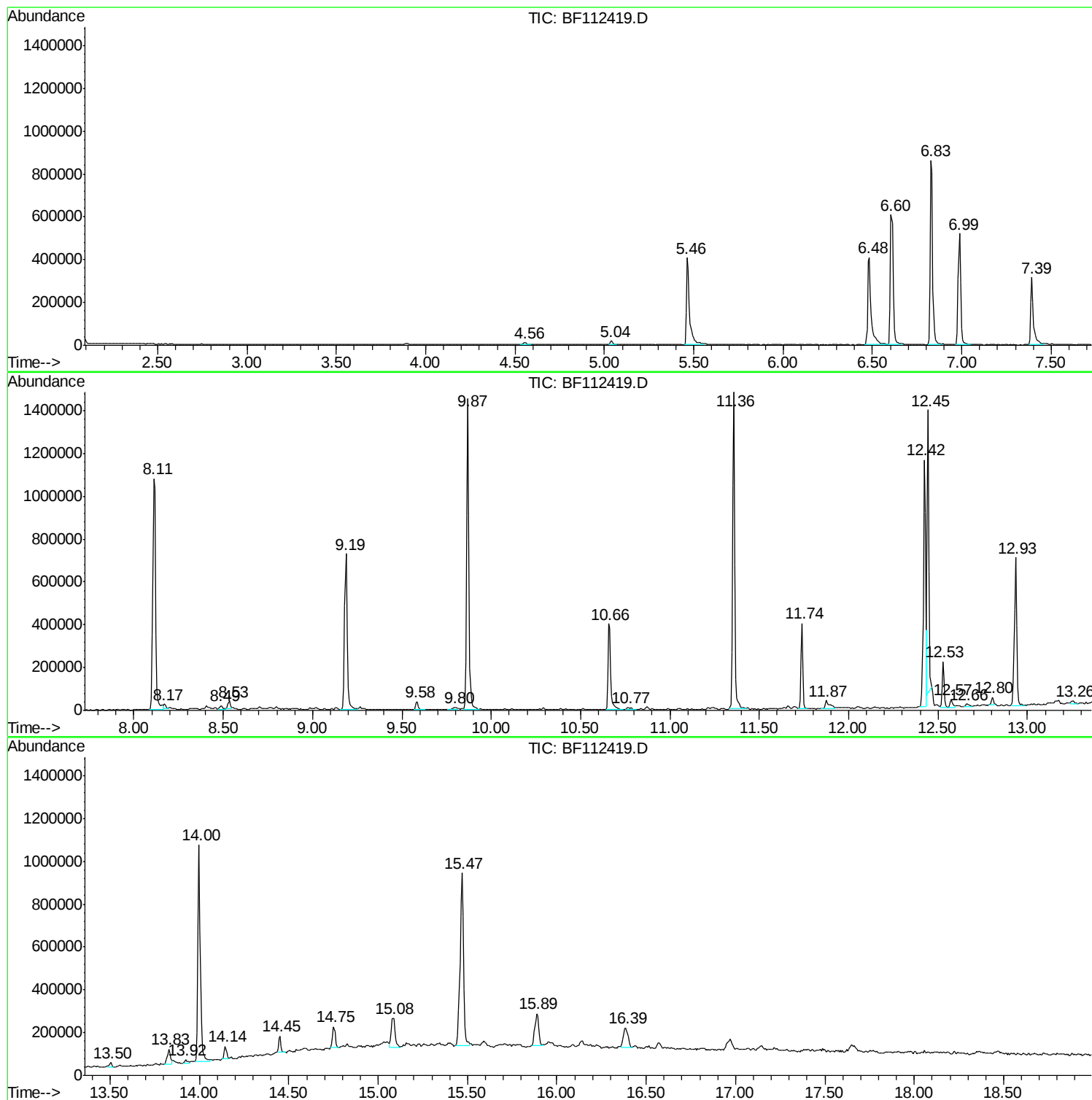
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF012519.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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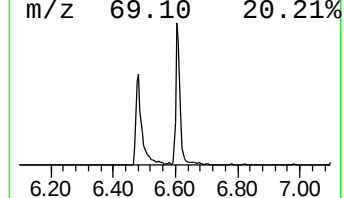
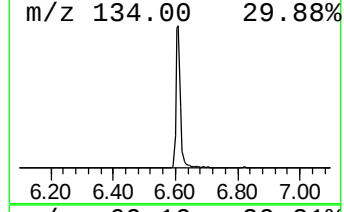
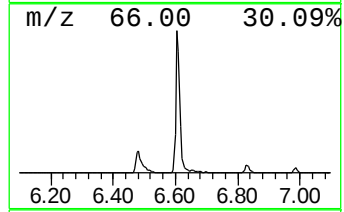
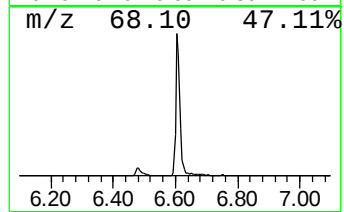
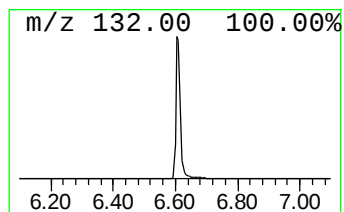
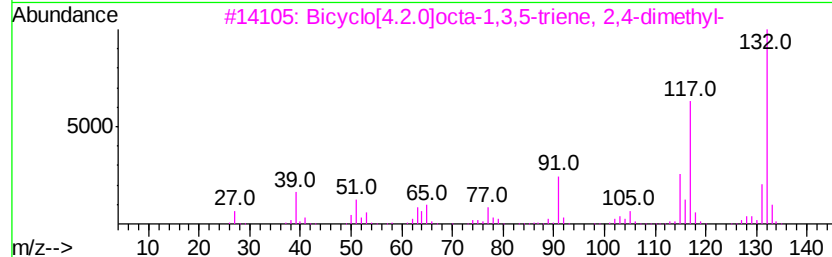
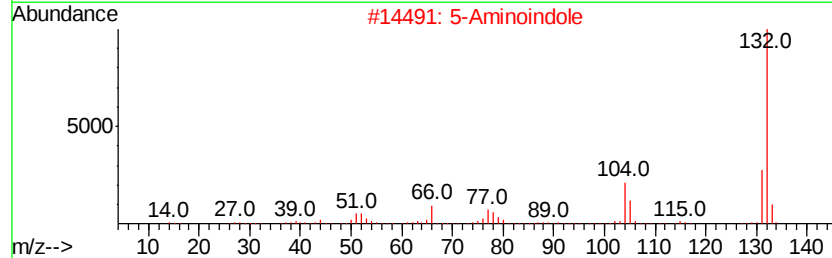
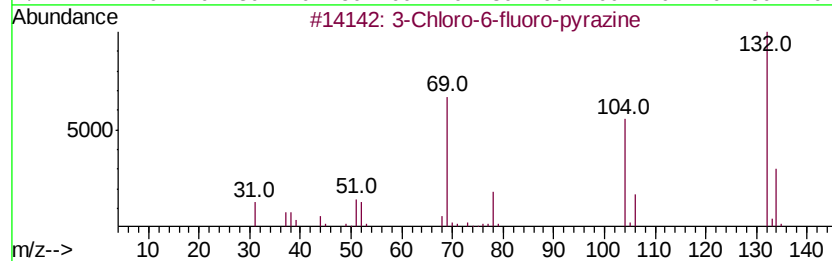
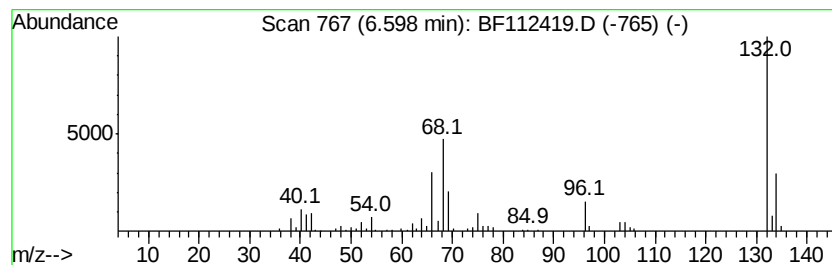
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF012519.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown6.60 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.60	14.50 ng	600717	1,4-Dichlorobenzene-d4	6.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
2		5-Aminoindole	132	C8H8N2	005192-03-0	12
3		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
4		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9
5		4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9



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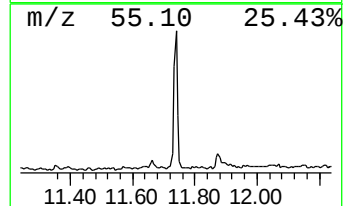
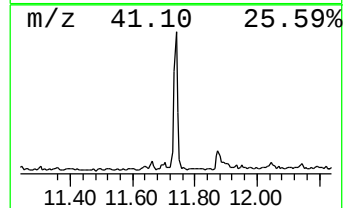
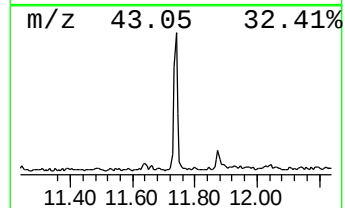
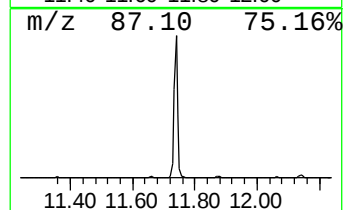
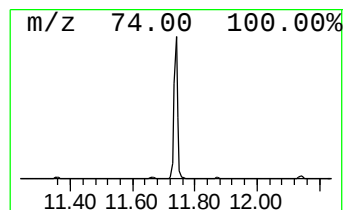
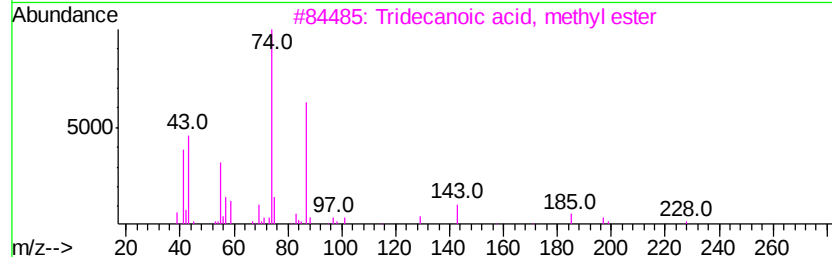
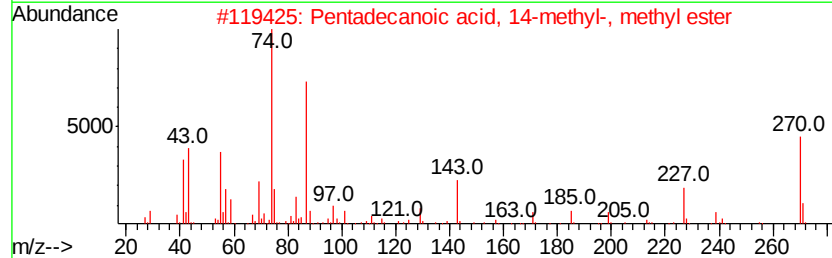
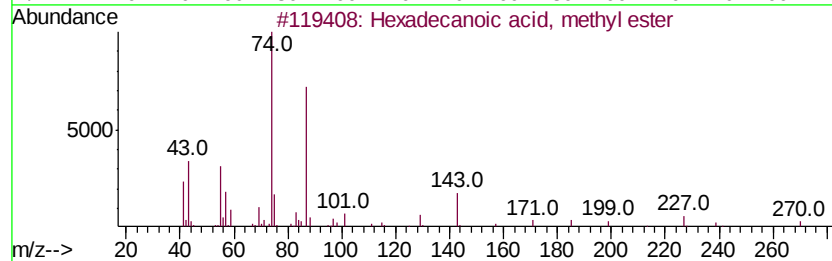
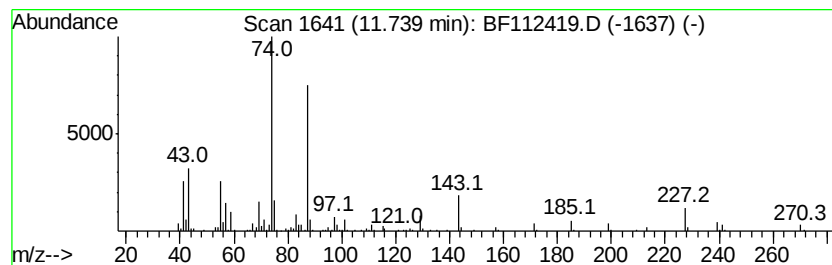
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 3 Hexadecanoic acid, methyl e... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.74	4.86 ng	322834	Phenanthrene-d10	11.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	98
2		Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	97
3		Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	93
4		Decanoic acid, methyl ester	186	C11H22O2	000110-42-9	87
5		Pentadecanoic acid, methyl ester	256	C16H32O2	007132-64-1	86



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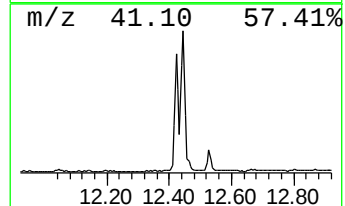
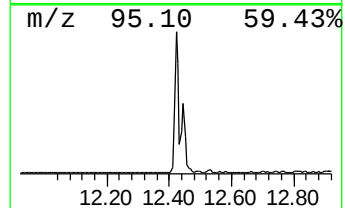
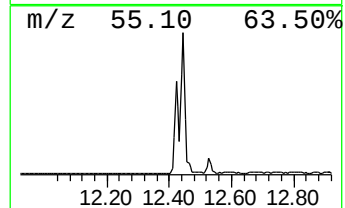
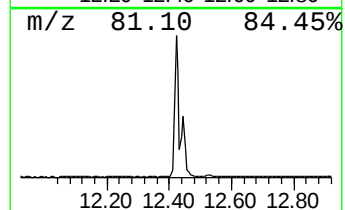
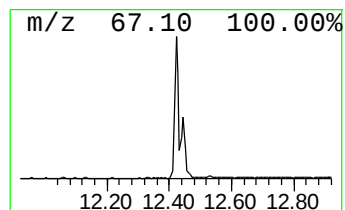
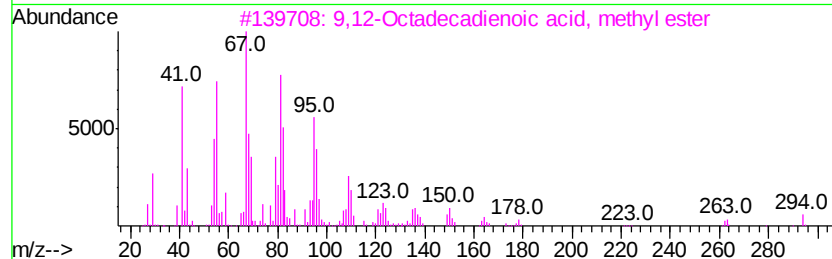
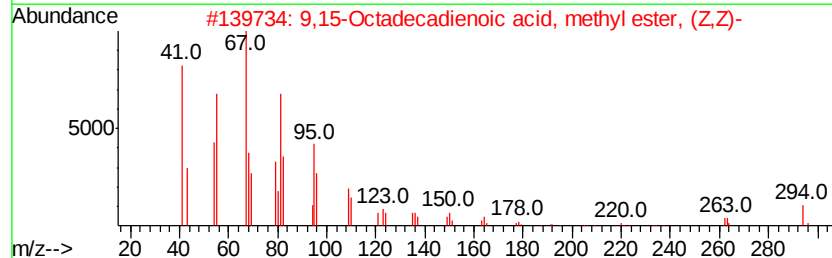
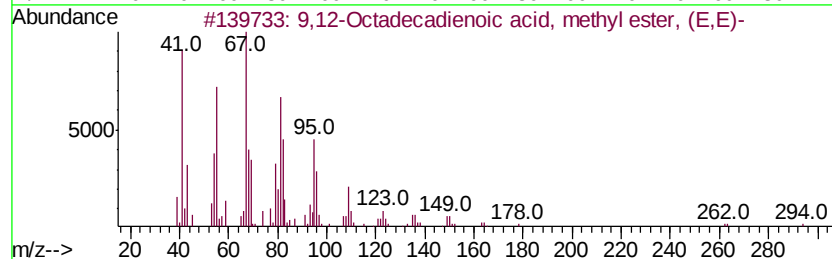
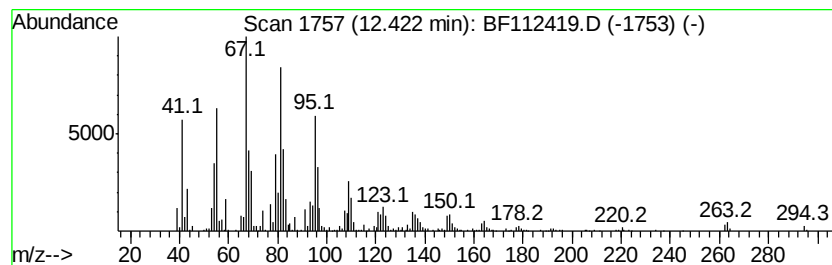
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 9,12-Octadecadienoic acid, ... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.		
12.42	15.36 ng	1020150	Phenanthrene-d10	11.36		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,12-Octadecadienoic acid, methy...	294	C19H34O2	002566-97-4	99	
2	9,15-Octadecadienoic acid, methy...	294	C19H34O2	017309-05-6	99	
3	9,12-Octadecadienoic acid, methy...	294	C19H34O2	002462-85-3	99	
4	10,13-Octadecadienoic acid, meth...	294	C19H34O2	056554-62-2	99	
5	9,12-Octadecadienoic acid (Z,Z)-...	294	C19H34O2	000112-63-0	99	



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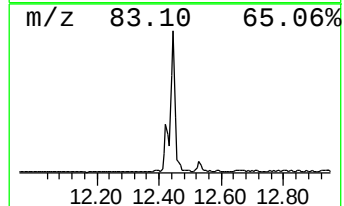
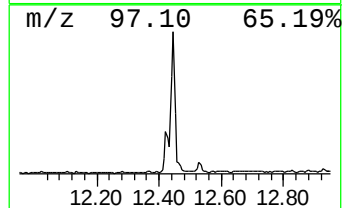
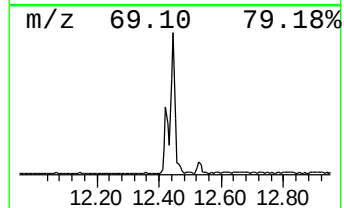
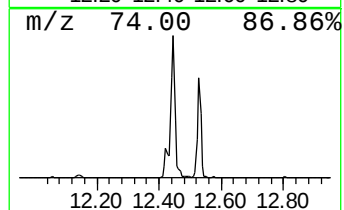
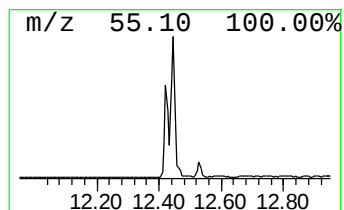
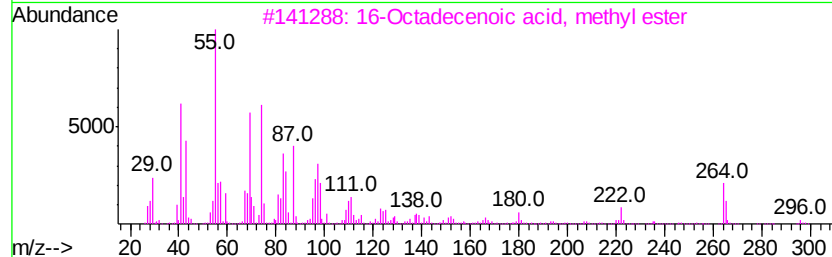
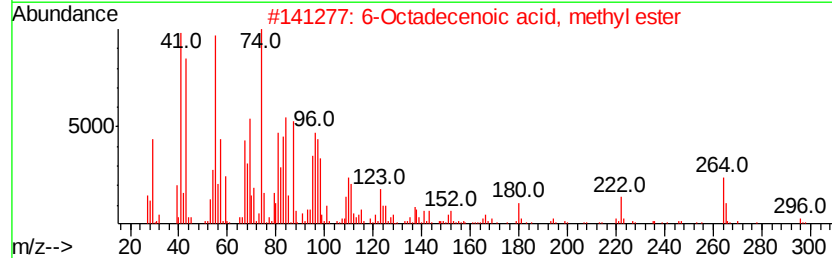
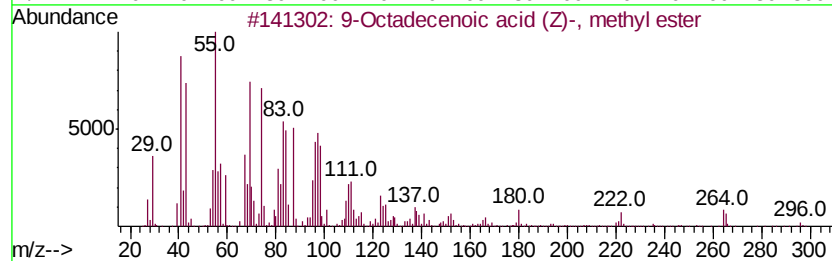
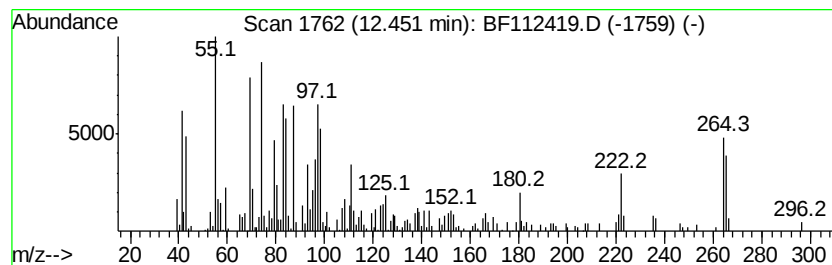
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TIC Integration Parameters: LSCINT.P

Peak Number 5 9-Octadecenoic acid (Z)-, m... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.45	14.98 ng	994683	Phenanthrene-d10	11.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9-Octadecenoic acid (Z)-, methyl...	296	C19H36O2	000112-62-9	99
2		6-Octadecenoic acid, methyl ester	296	C19H36O2	052355-31-4	99
3		16-Octadecenoic acid, methyl ester	296	C19H36O2	056554-49-5	99
4		9-Octadecenoic acid, methyl este...	296	C19H36O2	001937-62-8	99
5		10-Octadecenoic acid, methyl ester	296	C19H36O2	013481-95-3	99



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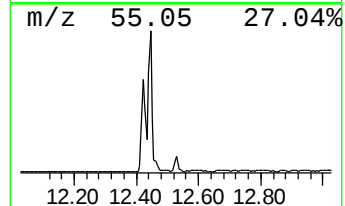
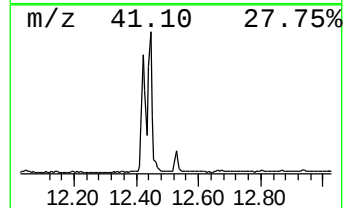
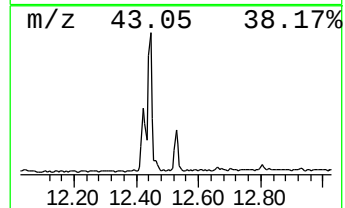
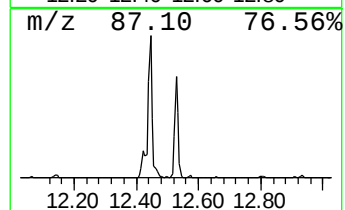
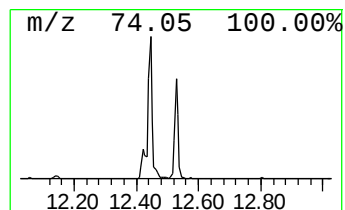
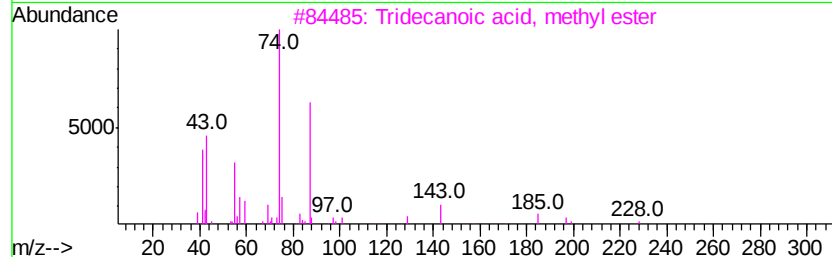
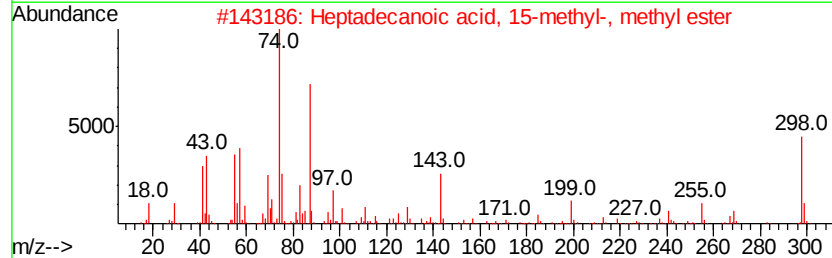
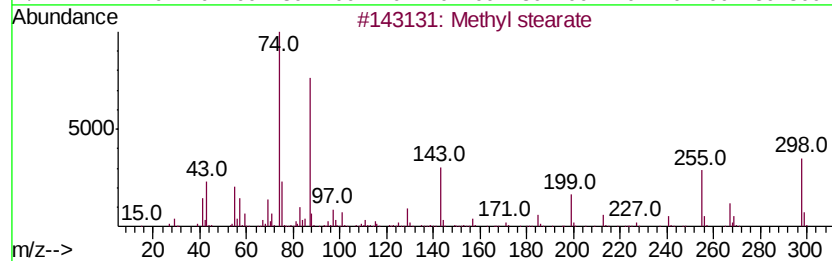
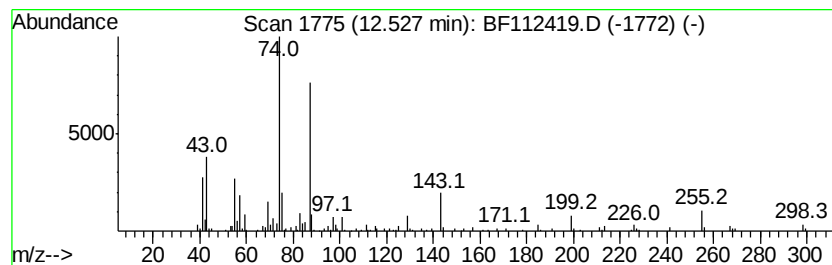
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Methyl stearate Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.53	2.62 ng	173710	Phenanthrene-d10	11.36
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Methyl stearate	298 C19H38O2	000112-61-8 98
2		Heptadecanoic acid, 15-methyl-, ...	298 C19H38O2	054833-55-5 93
3		Tridecanoic acid, methyl ester	228 C14H28O2	001731-88-0 93
4		Heptadecanoic acid, 16-methyl-, ...	298 C19H38O2	005129-61-3 90
5		Methyl tetradecanoate	242 C15H30O2	000124-10-7 83



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF020619\
Data File : BF112419.D
Acq On : 6 Feb 2019 14:52
Operator : JU/SJ
Sample : K1338-03 5X
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
AU-06-020519-B

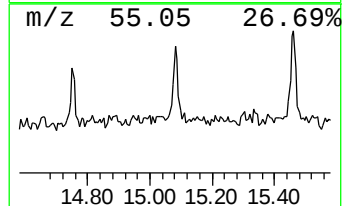
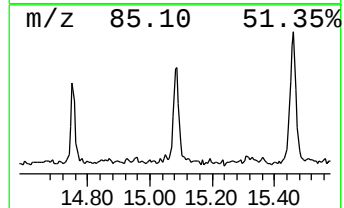
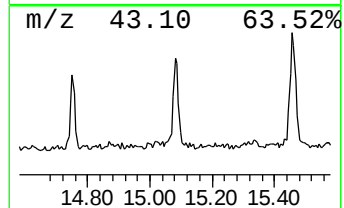
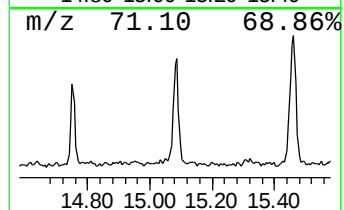
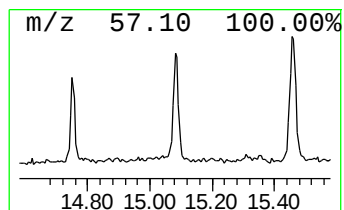
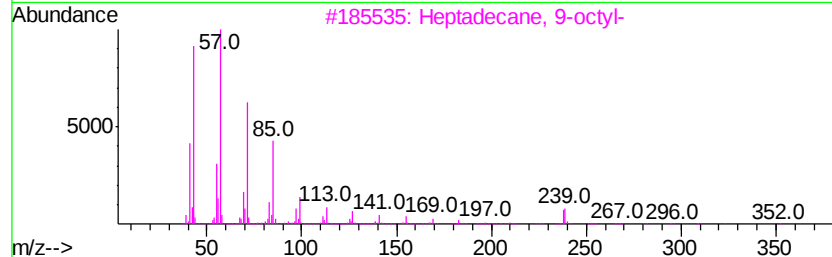
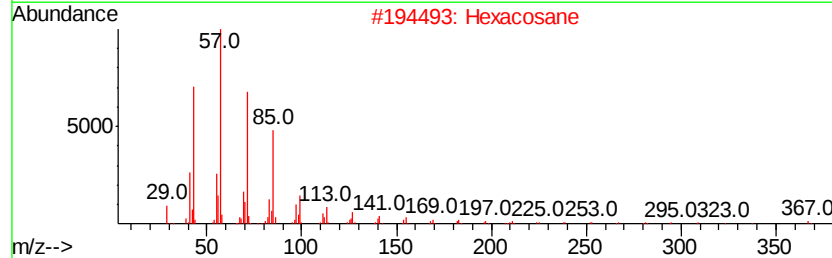
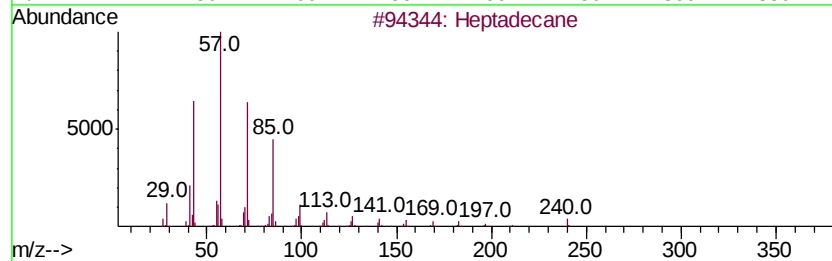
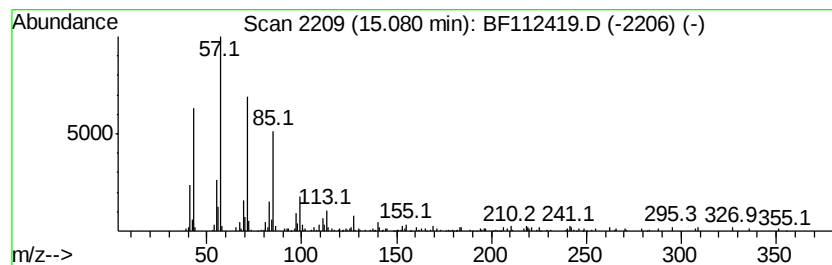
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Heptadecane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.08	3.33 ng	181780	Perylene-d12	15.47

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane	240	C17H36	000629-78-7	91
2		Hexacosane	366	C26H54	000630-01-3	90
3		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	87
4		2-methyloctacosane	408	C29H60	1000376-72-8	87
5		Heneicosane	296	C21H44	000629-94-7	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF020619\
Data File : BF112419.D
Acq On : 6 Feb 2019 14:52
Operator : JU/SJ
Sample : K1338-03 5X
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
AU-06-020519-B

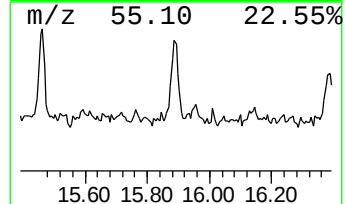
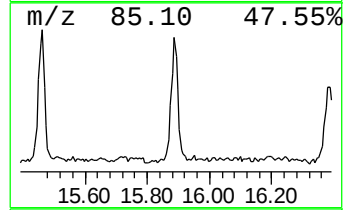
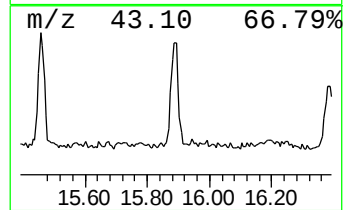
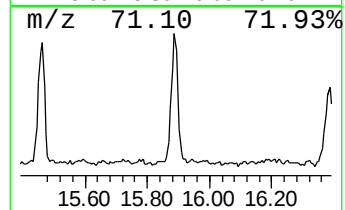
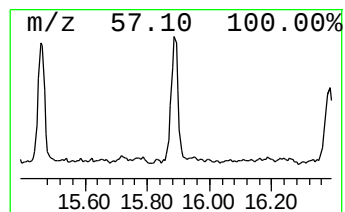
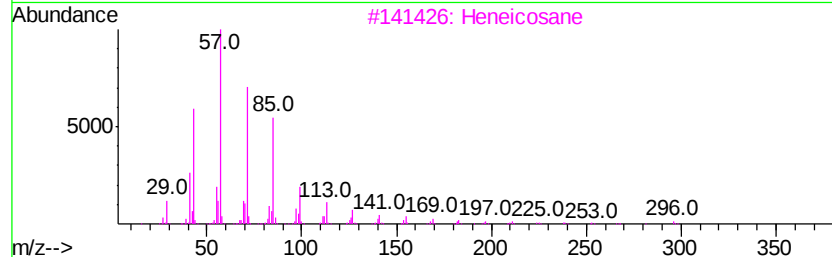
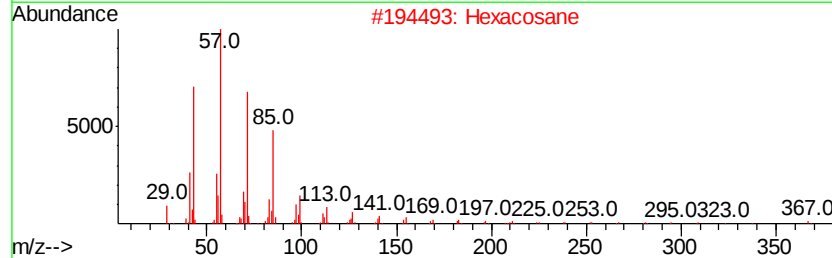
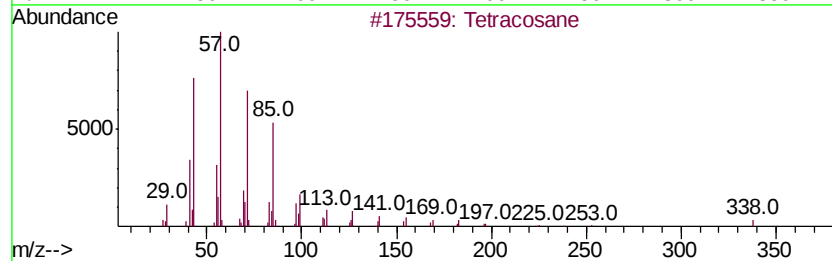
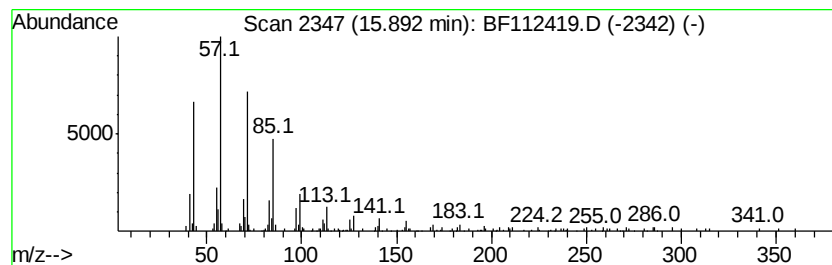
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Tetracosane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.89	3.98 ng	217379	Perylene-d12	15.47

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetracosane	338	C24H50	000646-31-1	91
2		Hexacosane	366	C26H54	000630-01-3	91
3		Heneicosane	296	C21H44	000629-94-7	91
4		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
5		Docosane, 7-hexyl-	394	C28H58	055373-86-9	91



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Sample : K1338-03 5X
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
AU-06-020519-B

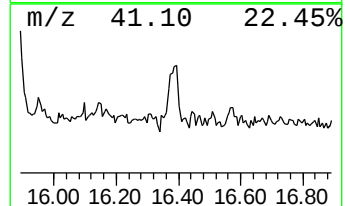
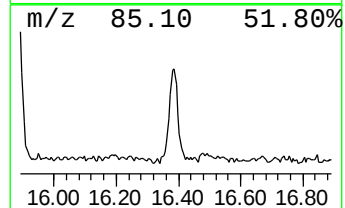
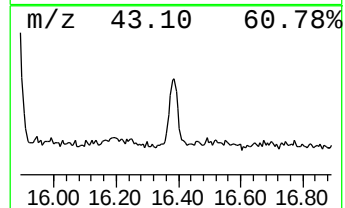
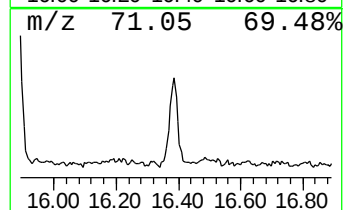
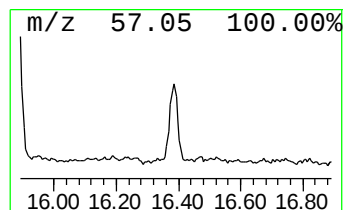
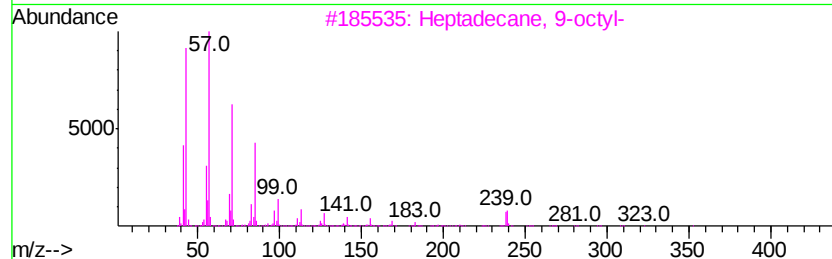
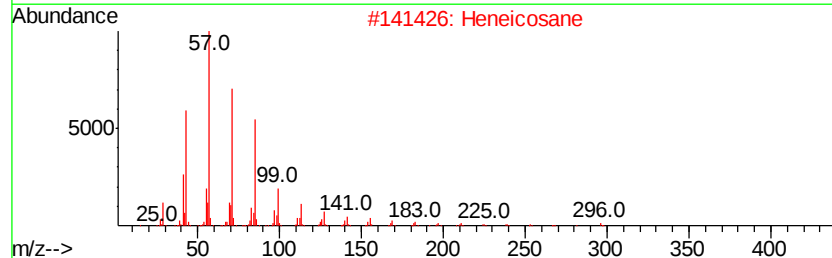
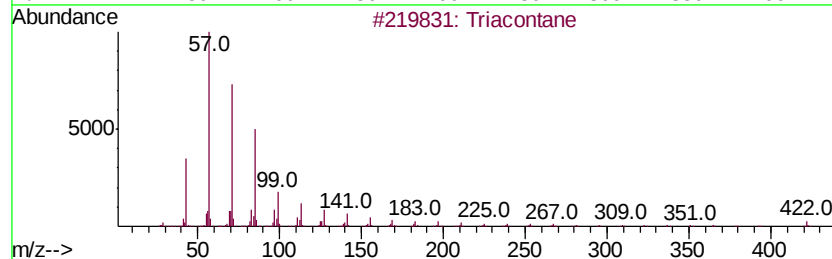
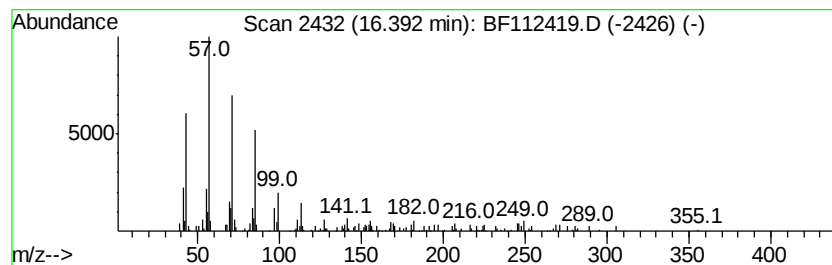
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Triacontane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.39	3.06 ng	167261	Perylene-d12	15.47

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Triacontane	422	C30H62	000638-68-6	87
2		Heneicosane	296	C21H44	000629-94-7	87
3		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	87
4		2-methyloctacosane	408	C29H60	1000376-72-8	87
5		Pentacosane	352	C25H52	000629-99-2	87



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Data File : BF112419.D
Acq On : 6 Feb 2019 14:52
Operator : JU/SJ
Sample : K1338-03 5X
Misc :
ALS Vial : 5 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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Hexadecanoic acid...	11.74	4.9	ng	322834	4	11.36	1328320	20.0
9,12-Octadecadien...	12.42	15.4	ng	1020150	4	11.36	1328320	20.0
9-Octadecenoic ac...	12.45	15.0	ng	994683	4	11.36	1328320	20.0
Methyl stearate	12.53	2.6	ng	173710	4	11.36	1328320	20.0
Heptadecane	15.08	3.3	ng	181780	6	15.47	1092540	20.0
Tetracosane	15.89	4.0	ng	217379	6	15.47	1092540	20.0
Triacontane	16.39	3.1	ng	167261	6	15.47	1092540	20.0