

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF101618\
 Data File : BF109904.D
 Acq On : 16 Oct 2018 14:46
 Operator : JU/SJ
 Sample : J5526-10
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SUP-4B-C-101518-1

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-BF101518.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	2.340	37	43	55	rVB	1002587	1386164	18.76%	2.182%
2	3.598	254	257	274	rBV	126818	245871	3.33%	0.387%
3	5.234	530	535	544	rVB	254380	348538	4.72%	0.549%
4	5.610	593	599	602	rBV	6484809	7033073	95.19%	11.072%
5	6.604	762	768	772	rBV	6248689	7388355	100.00%	11.632%
6	6.751	788	793	796	rBV	6923903	7148991	96.76%	11.255%
7	6.981	829	832	836	rVB	2147151	1650357	22.34%	2.598%
8	7.139	855	859	862	rVB	6444053	5600282	75.80%	8.817%
9	7.545	922	928	931	rBV	4573681	4640063	62.80%	7.305%
10	8.263	1046	1050	1053	rBV	2453807	2016147	27.29%	3.174%
11	9.339	1228	1233	1236	rBV	7115175	6703805	90.73%	10.554%
12	9.727	1295	1299	1302	rBV	1035639	799541	10.82%	1.259%
13	10.022	1345	1349	1353	rBV	2230550	1865271	25.25%	2.937%
14	10.810	1479	1483	1487	rBV	3246643	3023703	40.93%	4.760%
15	11.516	1599	1603	1606	rBV	1574796	1458981	19.75%	2.297%
16	12.016	1684	1688	1691	rBV	78511	94727	1.28%	0.149%
17	12.539	1774	1777	1780	rBV	181701	149787	2.03%	0.236%
18	12.727	1806	1809	1815	rBV	97193	106418	1.44%	0.168%
19	12.804	1818	1822	1828	rBV2	98781	161195	2.18%	0.254%
20	12.957	1846	1848	1851	rVB	101847	78536	1.06%	0.124%
21	13.098	1868	1872	1875	rBV	5110649	4708252	63.73%	7.412%
22	13.974	2018	2021	2033	rVB2	485804	577675	7.82%	0.909%
23	14.157	2048	2052	2055	rBV	1983285	1767363	23.92%	2.782%
24	14.298	2073	2076	2079	rBV	168451	167201	2.26%	0.263%
25	14.609	2125	2129	2134	rVB	338508	310355	4.20%	0.489%
26	14.933	2180	2184	2193	rVB	339022	377214	5.11%	0.594%
27	15.286	2240	2244	2248	rBV	427328	460604	6.23%	0.725%
28	15.686	2307	2312	2324	rBV2	1675716	2342270	31.70%	3.687%
29	16.156	2388	2392	2398	rVV	314214	466283	6.31%	0.734%
30	16.215	2398	2402	2408	rVB4	68935	108992	1.48%	0.172%
31	16.703	2480	2485	2494	rVB2	165772	333485	4.51%	0.525%

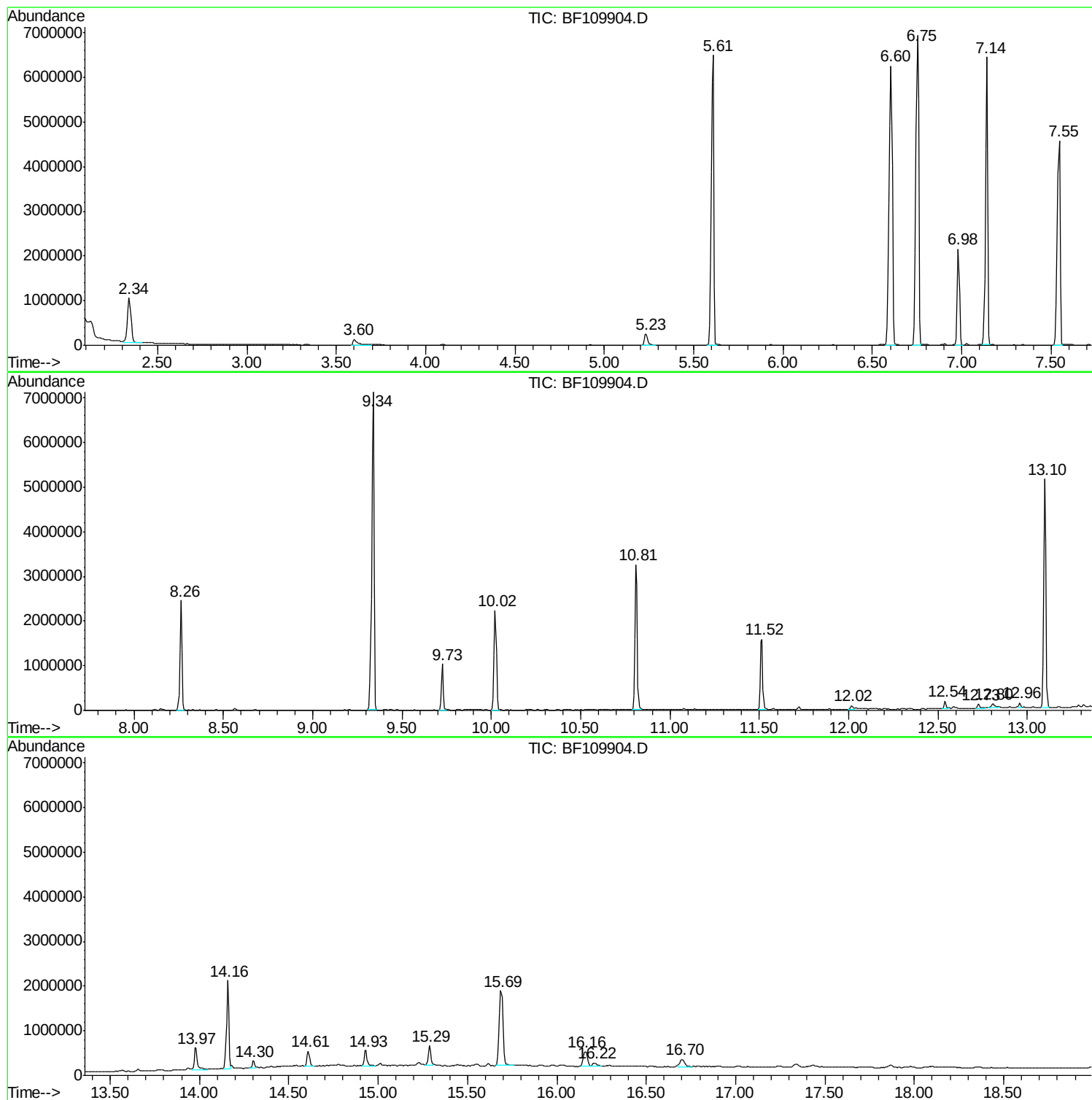
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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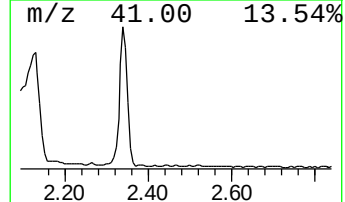
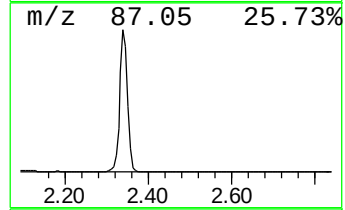
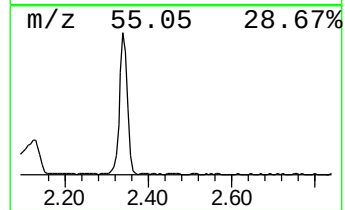
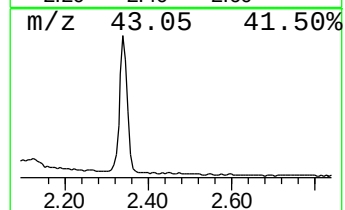
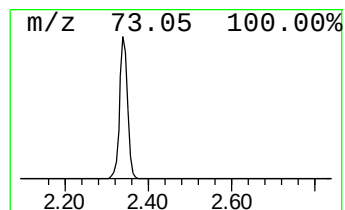
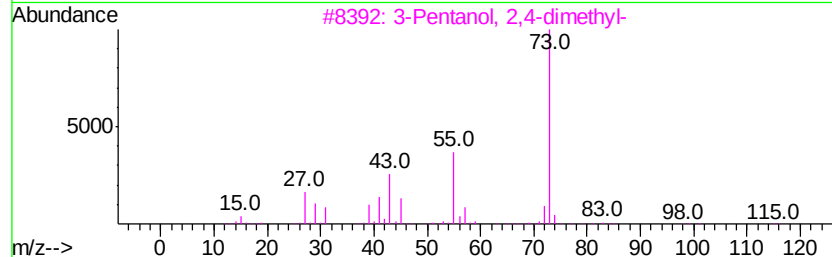
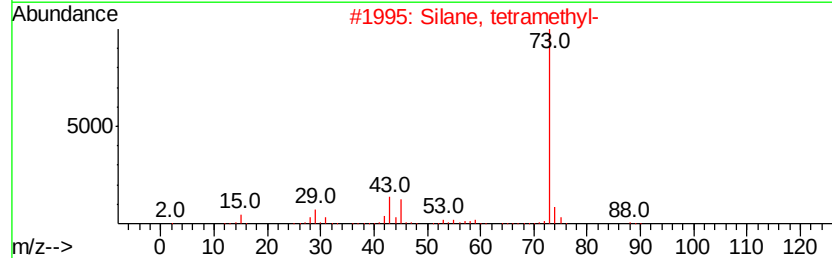
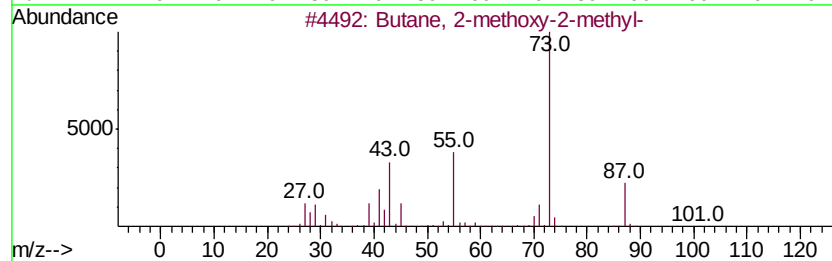
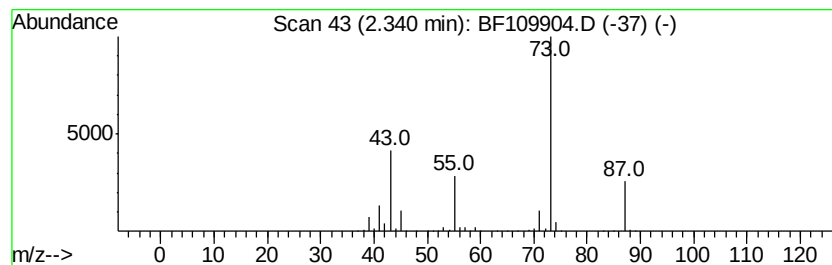
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.34	16.80 ng	1386160	1,4-Dichlorobenzene-d4	6.98

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2			Silane, tetramethyl-	88	C4H12Si	000075-76-3	38
3			3-Pentanol, 2,4-dimethyl-	116	C7H16O	000600-36-2	37
4			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	23
5			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12



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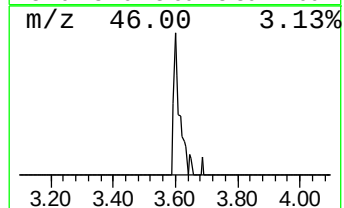
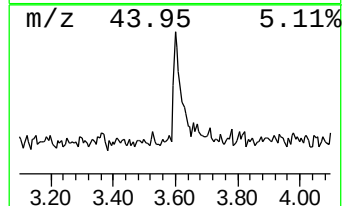
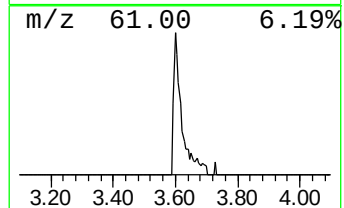
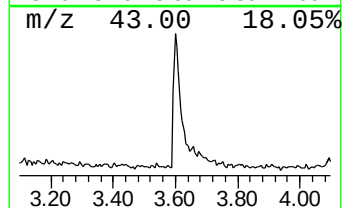
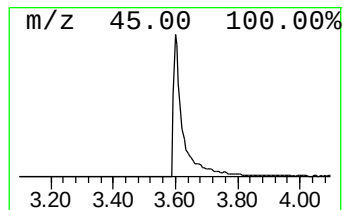
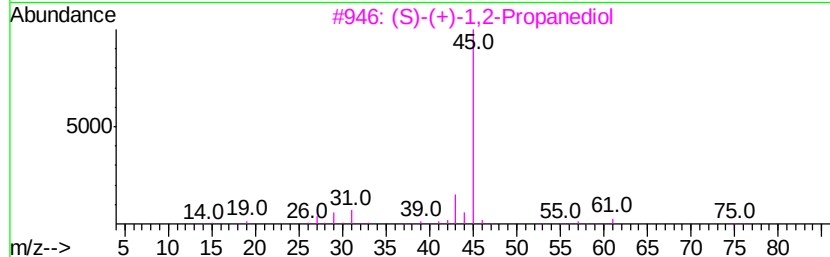
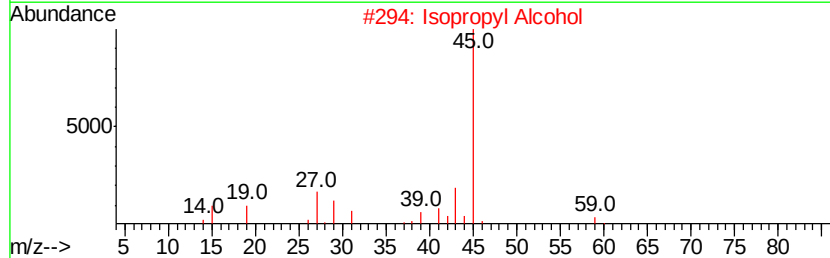
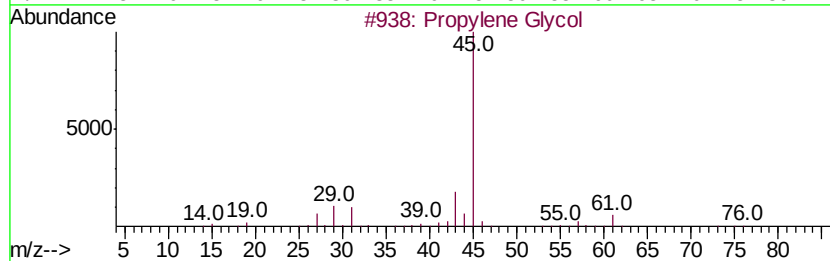
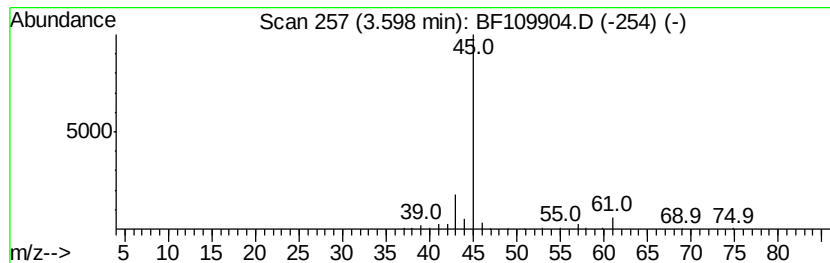
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 2 unknown3.60 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.60	2.98 ng	245871	1,4-Dichlorobenzene-d4	6.98

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propylene Glycol	76	C3H8O2	000057-55-6	43
2		Isopropyl Alcohol	60	C3H8O	000067-63-0	9
3		(S)-(+)-1,2-Propanediol	76	C3H8O2	004254-15-3	9
4		Propanoic acid, 2-hydroxy-, meth...	104	C4H8O3	002155-30-8	9
5		Formamide	45	CH3NO	000075-12-7	7



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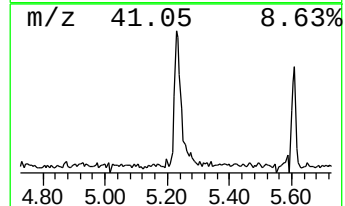
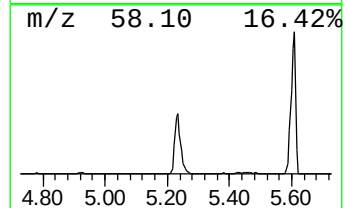
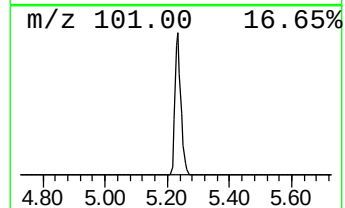
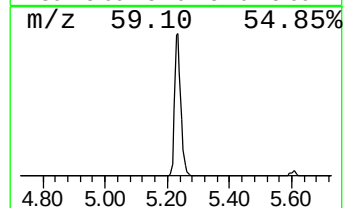
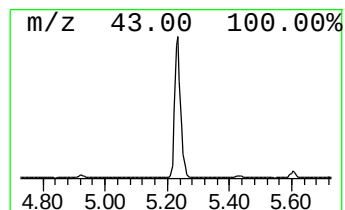
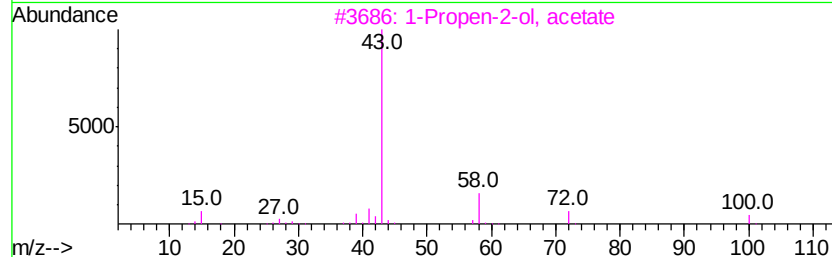
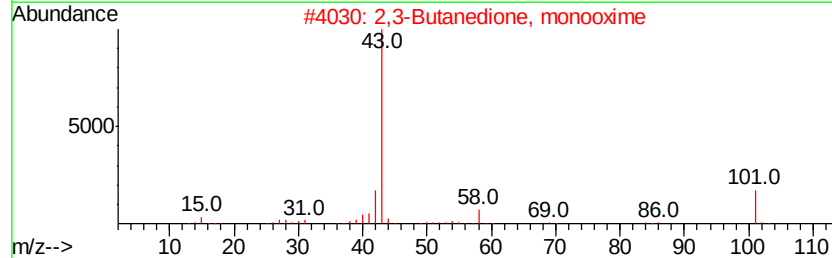
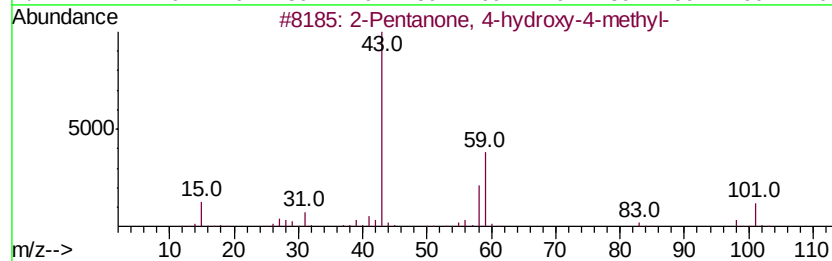
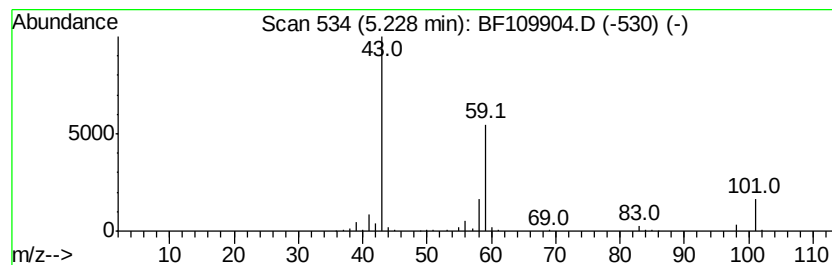
Instrument :
BNA_F
ClientSampleId :
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TIC Integration Parameters: LSCINT.P

Peak Number 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
5.23	4.22 ng	348538	1,4-Dichlorobenzene-d4	6.98	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16
3	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
4	5-Oxohexanenitrile	111	C6H9NO	010412-98-3	9
5	Ethanol, pentamethyl-	116	C7H16O	000594-83-2	9



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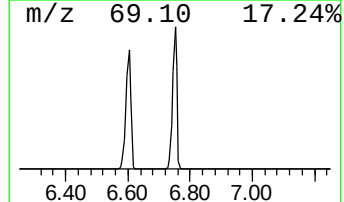
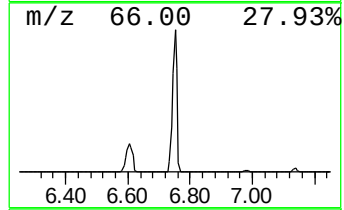
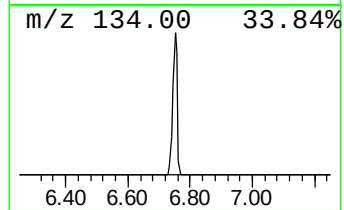
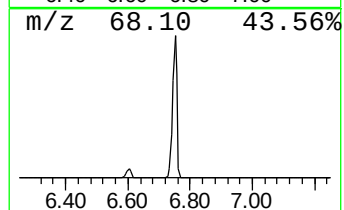
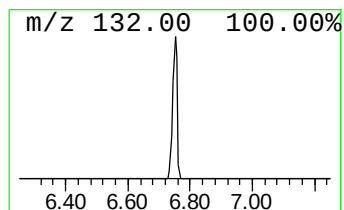
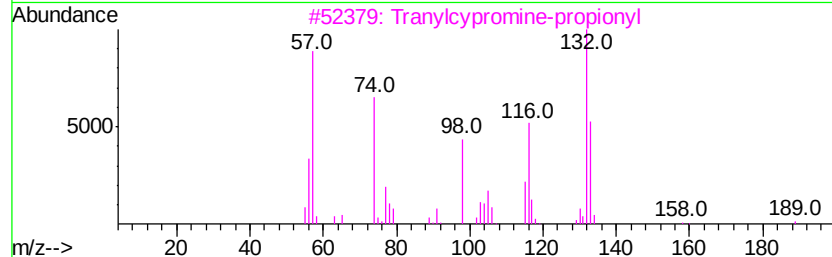
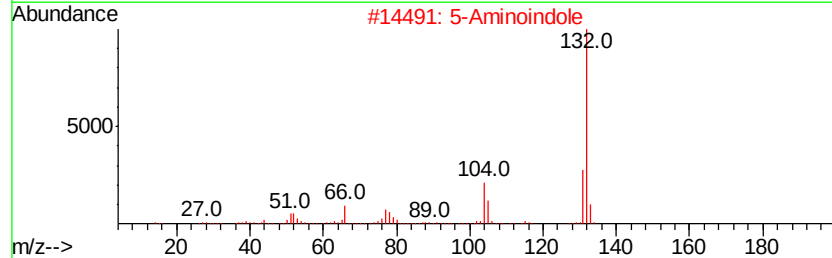
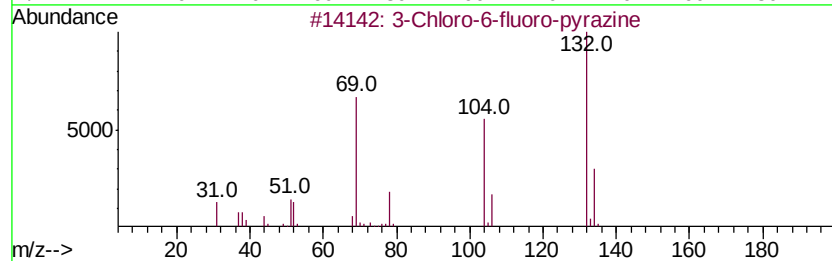
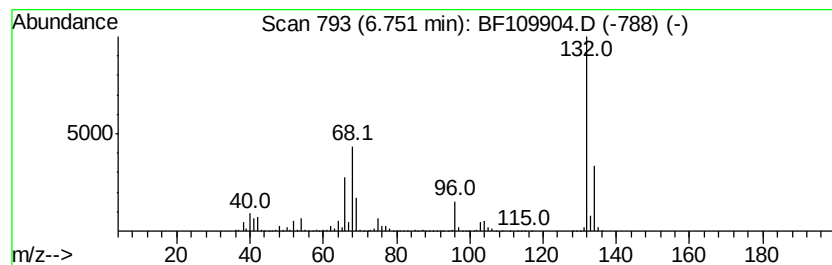
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 unknown6.75 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.75	86.64 ng	7148990	1,4-Dichlorobenzene-d4	6.98

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		5-Aminoindole	132	C8H8N2	005192-03-0	12
3		Tranvlcyproline-propionyl	189	C12H15NO	1000123-86-3	10
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9



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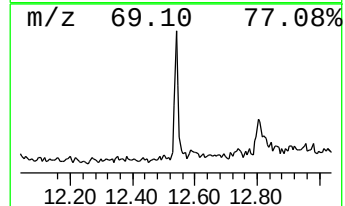
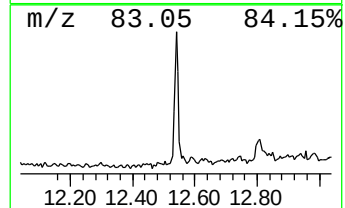
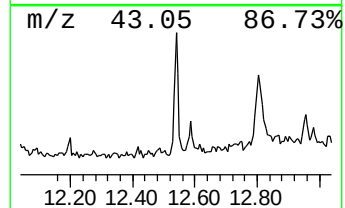
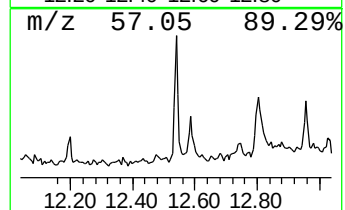
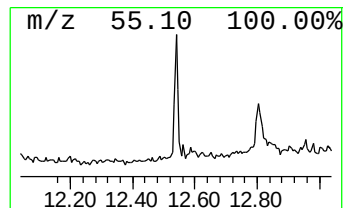
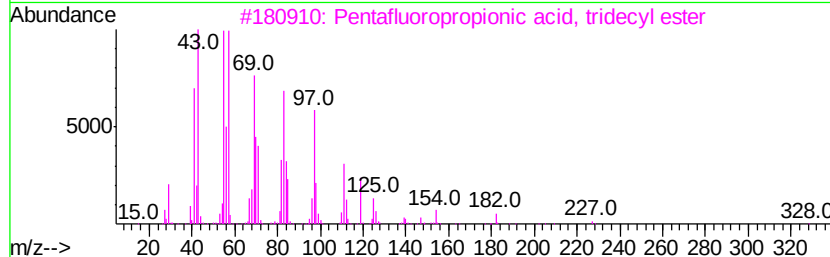
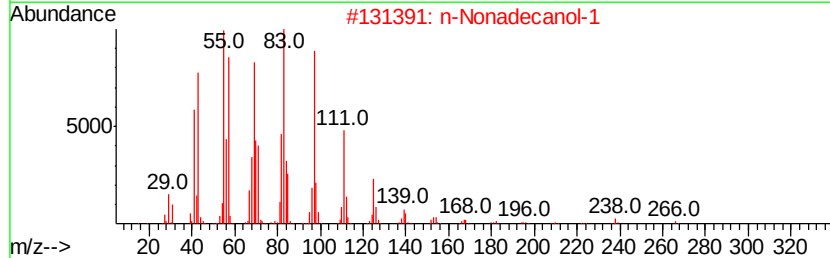
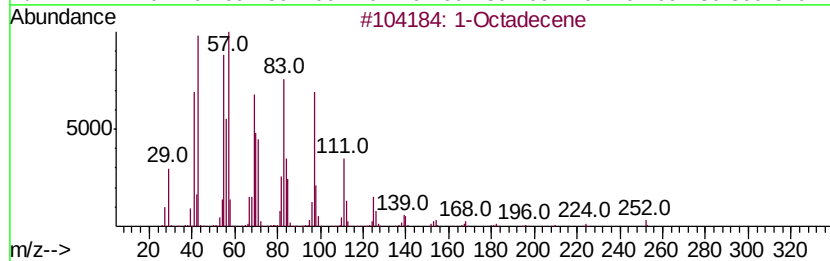
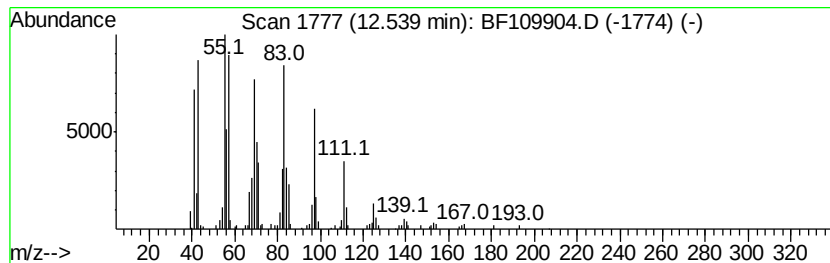
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 6 1-Octadecene Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.54	2.05 ng	149787	Phenanthrene-d10	11.52

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Octadecene	252	C18H36	000112-88-9	93
2		n-Nonadecanol-1	284	C19H40O	001454-84-8	91
3		Pentafluoropropionic acid, tridecyl ester	346	C16H27F5O2	959261-22-4	91
4		Pentafluoropropionic acid, dodecyl ester	332	C15H25F5O2	006222-04-4	91
5		1-Tetradecanol	214	C14H30O	000112-72-1	91



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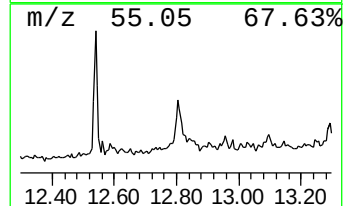
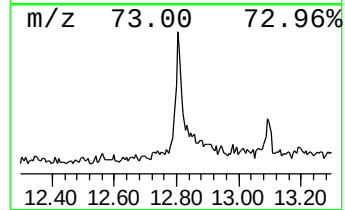
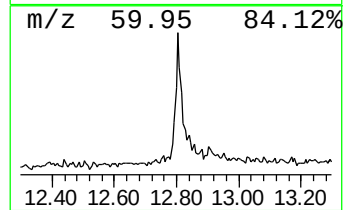
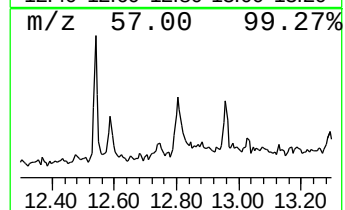
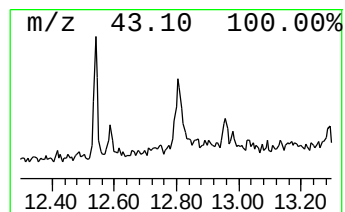
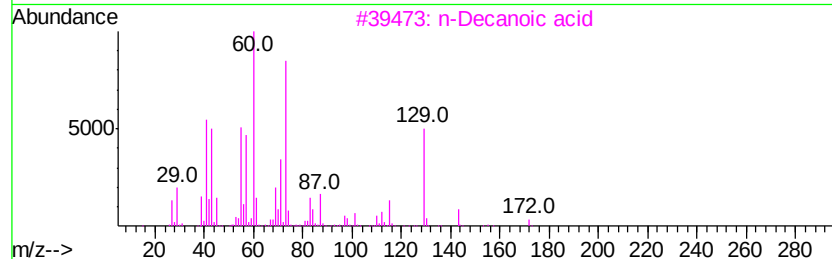
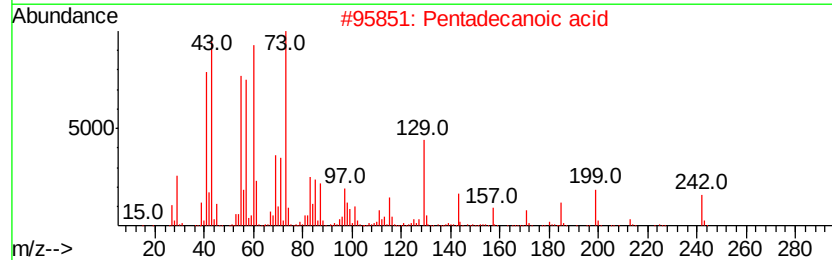
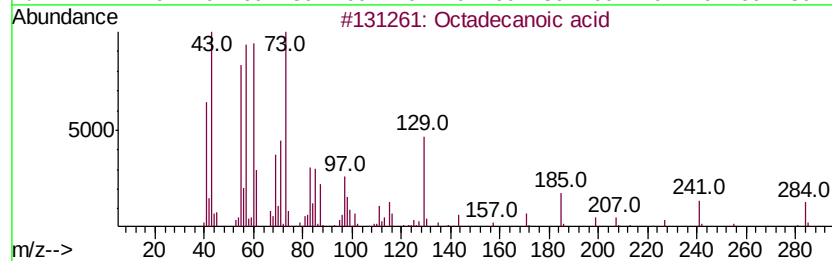
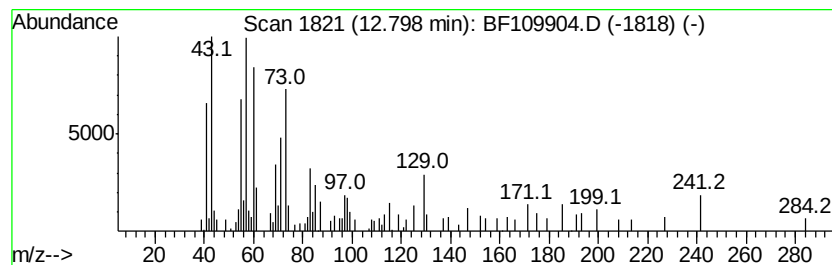
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Octadecanoic acid Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD		R.T.
12.80	2.21 ng	161195	Phenanthrene-d10		11.52
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	96
2	Pentadecanoic acid	242	C15H30O2	001002-84-2	80
3	n-Decanoic acid	172	C10H20O2	000334-48-5	76
4	Tridecanoic acid	214	C13H26O2	000638-53-9	64
5	Undecanoic acid	186	C11H22O2	000112-37-8	59



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101618\
Data File : BF109904.D
Acq On : 16 Oct 2018 14:46
Operator : JU/SJ
Sample : J5526-10
Misc :
ALS Vial : 6 Sample Multiplier: 1

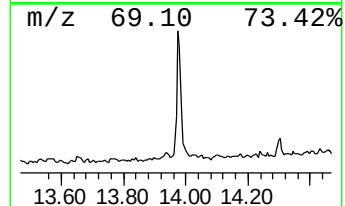
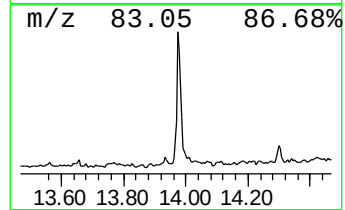
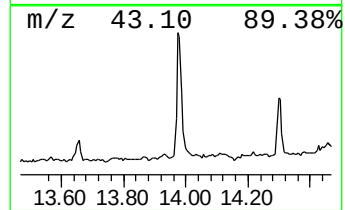
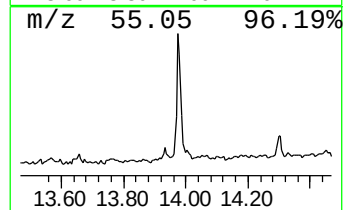
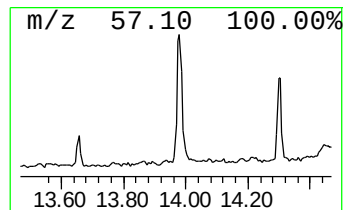
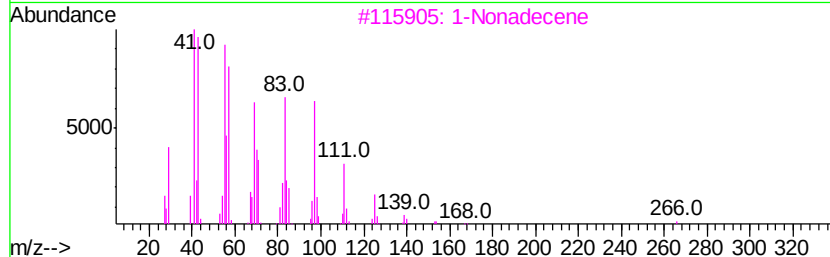
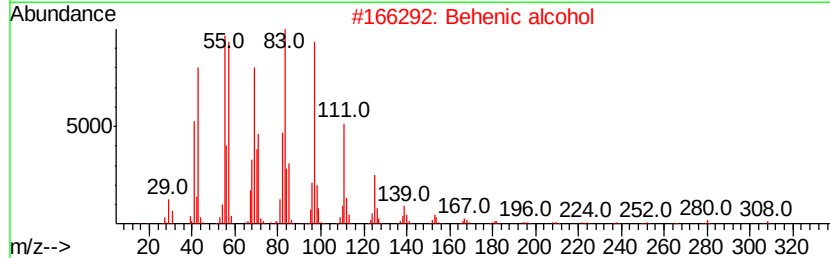
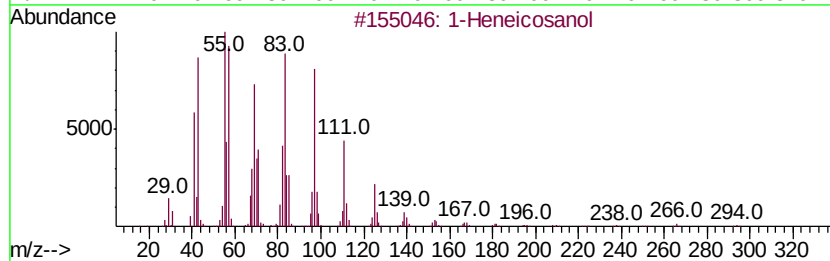
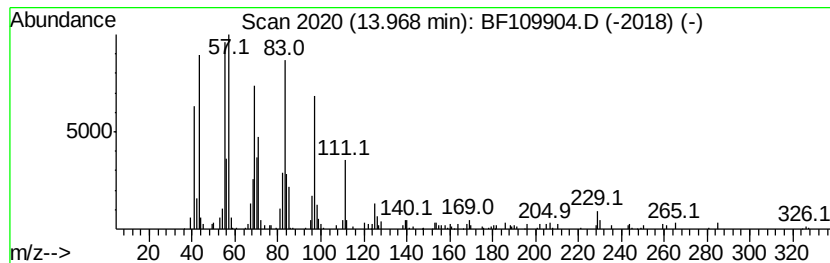
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 1-Heneicosanol Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.97	6.54 ng	577675	Chrysene-d12	14.16	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heneicosanol	312	C21H44O	015594-90-8	95
2	Behenic alcohol	326	C22H46O	000661-19-8	91
3	1-Nonadecene	266	C19H38	018435-45-5	91
4	3-Eicosene, (E)-	280	C20H40	074685-33-9	91
5	n-Tetracosanol-1	354	C24H50O	000506-51-4	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101618\
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Operator : JU/SJ
Sample : J5526-10
Misc :
ALS Vial : 6 Sample Multiplier: 1

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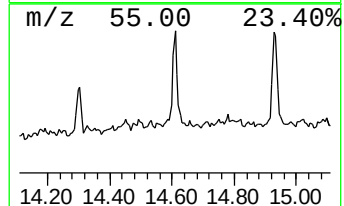
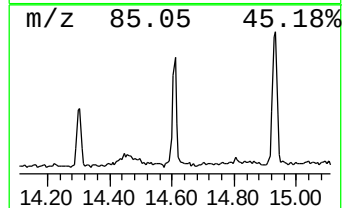
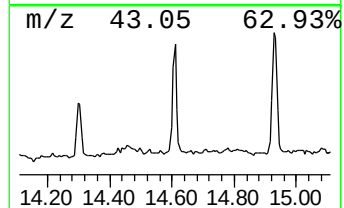
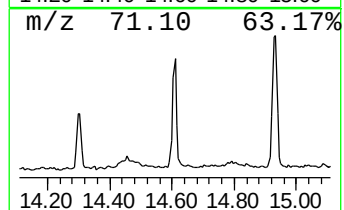
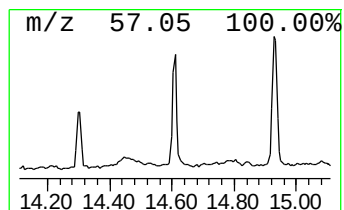
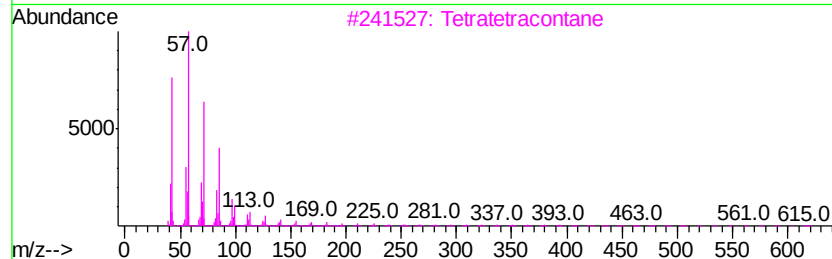
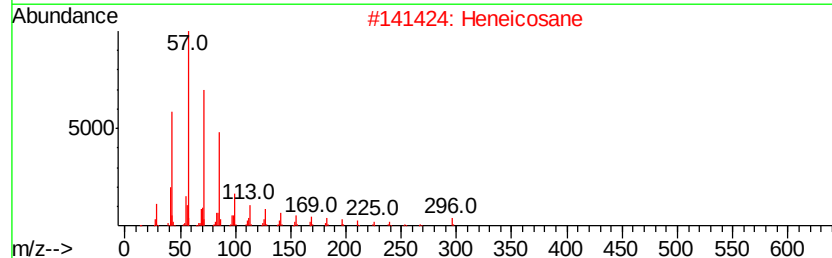
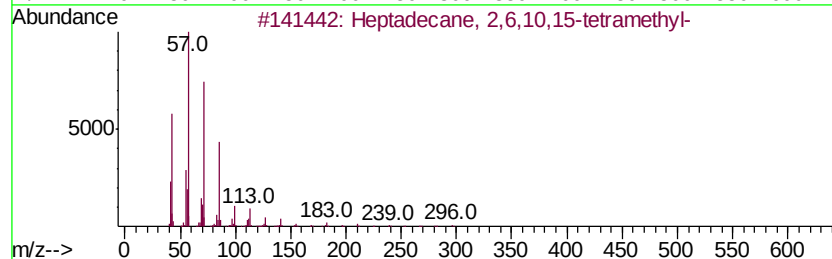
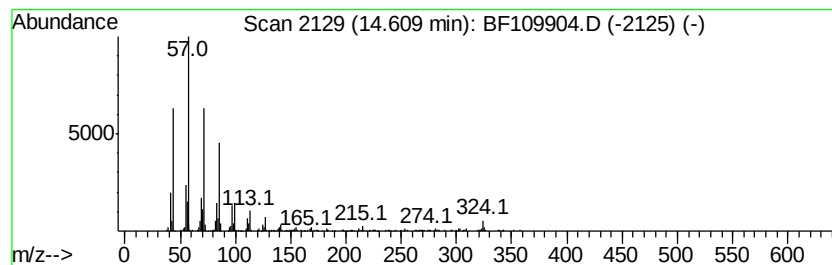
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Heptadecane, 2,6,10,15-tetr... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.61	3.51 ng	310355	Chrysene-d12	14.16

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	95
2			Heneicosane	296	C21H44	000629-94-7	91
3			Tetratetracontane	619	C44H90	007098-22-8	90
4			Octacosane	394	C28H58	000630-02-4	90
5			Tricosane	324	C23H48	000638-67-5	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101618\
Data File : BF109904.D
Acq On : 16 Oct 2018 14:46
Operator : JU/SJ
Sample : J5526-10
Misc :
ALS Vial : 6 Sample Multiplier: 1

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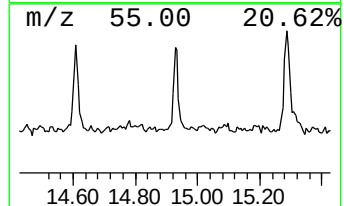
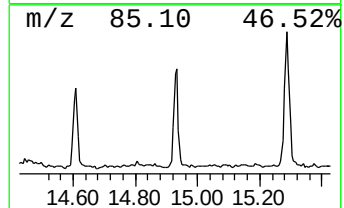
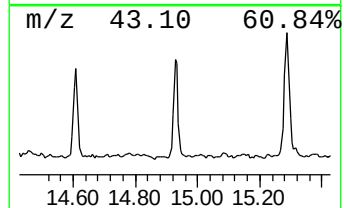
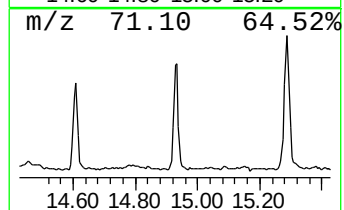
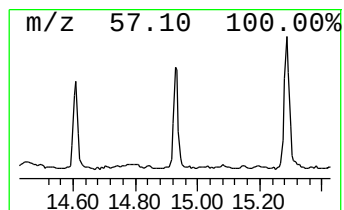
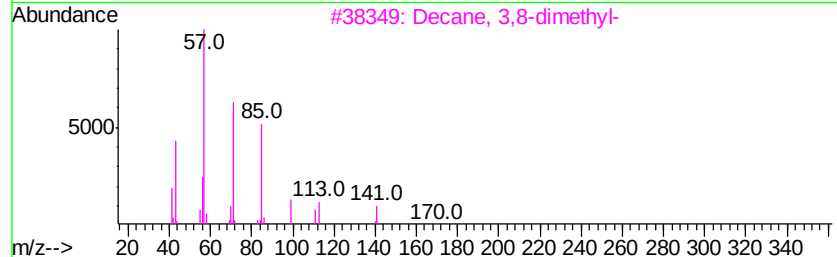
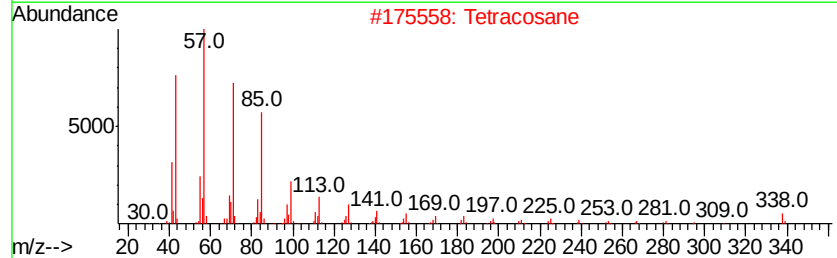
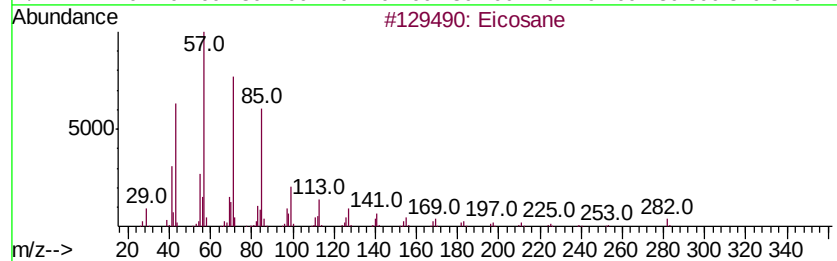
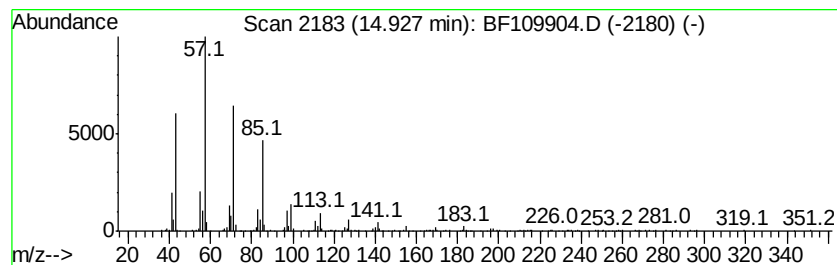
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Eicosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.93	3.22 ng	377214	Perylene-d12	15.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	91
2		Tetracosane	338	C24H50	000646-31-1	90
3		Decane, 3,8-dimethyl-	170	C12H26	017312-55-9	83
4		Hexadecane	226	C16H34	000544-76-3	83
5		Hexacosane	366	C26H54	000630-01-3	76



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101618\
Data File : BF109904.D
Acq On : 16 Oct 2018 14:46
Operator : JU/SJ
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ALS Vial : 6 Sample Multiplier: 1

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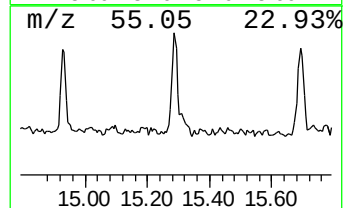
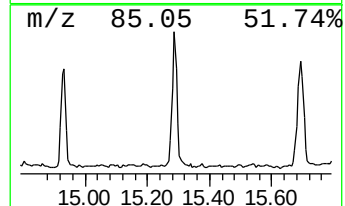
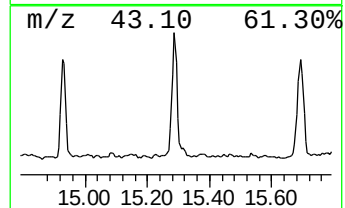
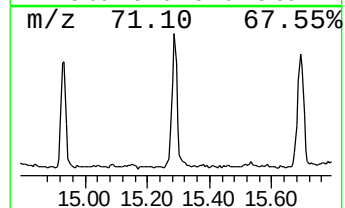
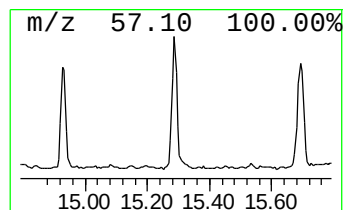
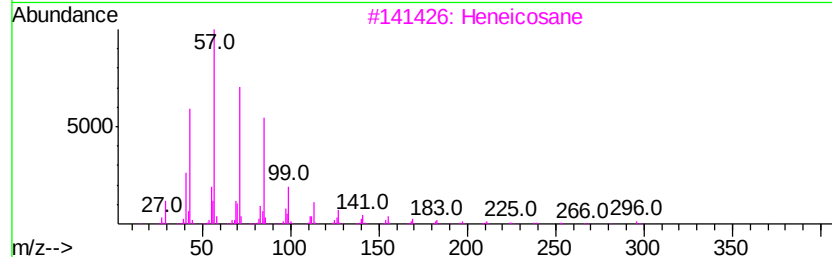
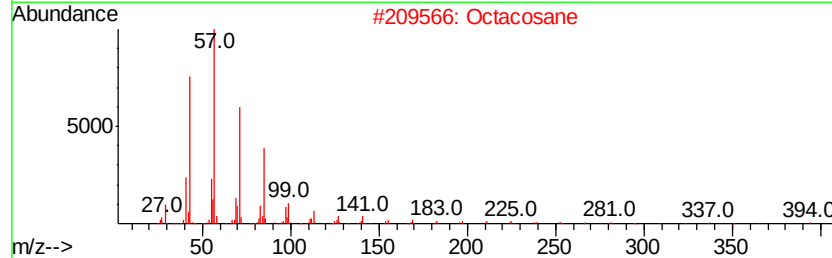
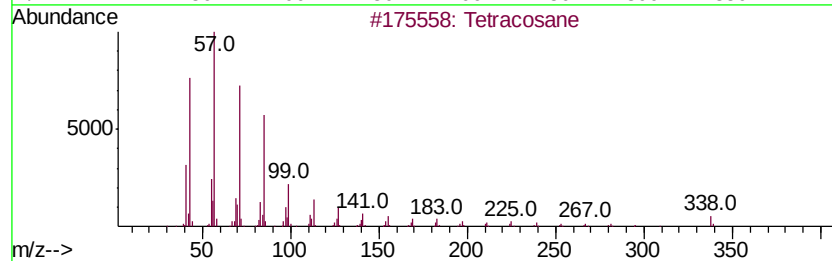
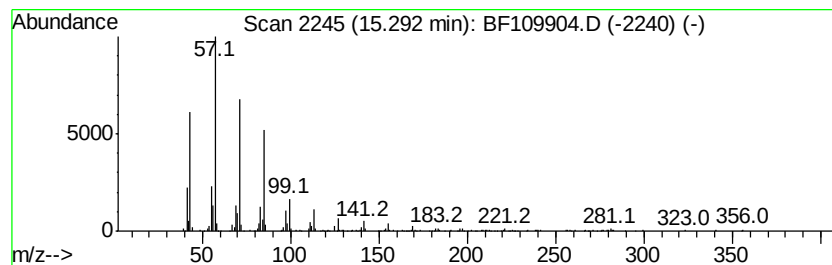
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Tetracosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.29	3.93 ng	460604	Perylene-d12	15.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetracosane	338	C24H50	000646-31-1	97
2		Octacosane	394	C28H58	000630-02-4	91
3		Heneicosane	296	C21H44	000629-94-7	91
4		Octadecane, 1-iodo-	380	C18H37I	000629-93-6	87
5		Hexacosane	366	C26H54	000630-01-3	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101618\
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Acq On : 16 Oct 2018 14:46
Operator : JU/SJ
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ALS Vial : 6 Sample Multiplier: 1

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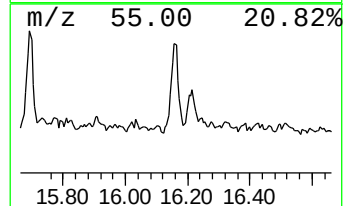
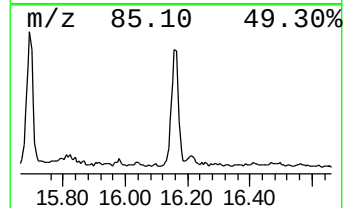
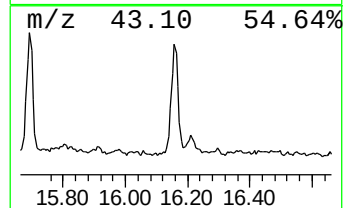
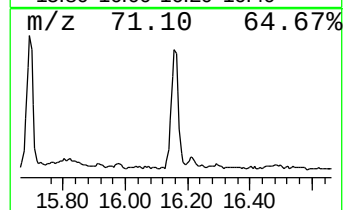
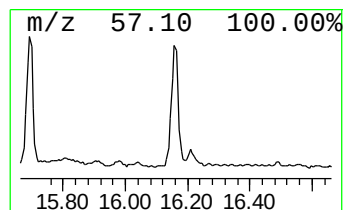
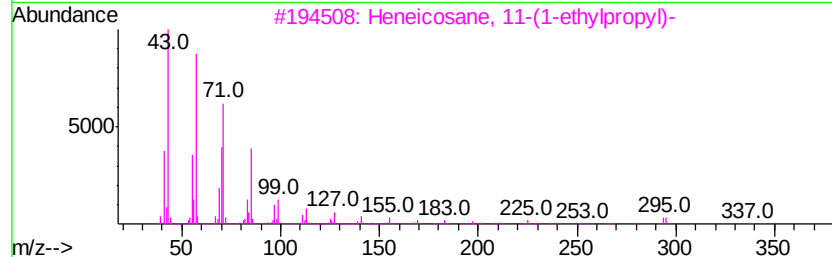
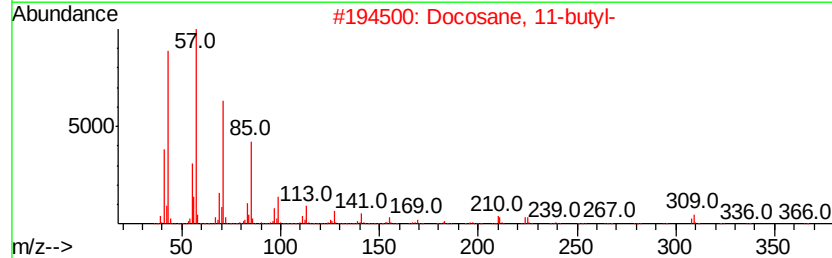
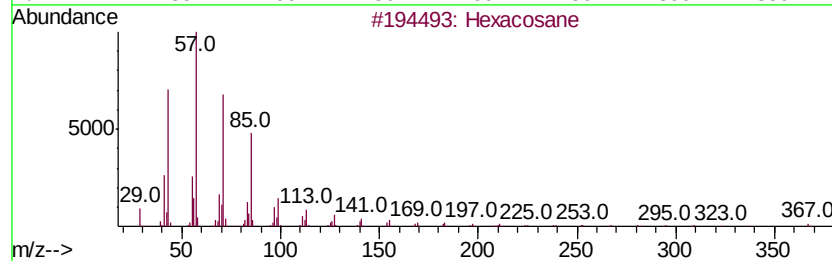
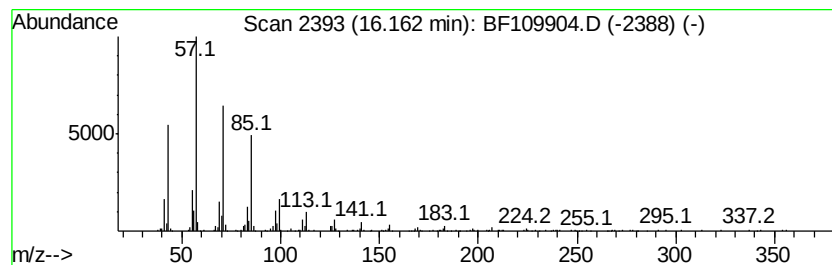
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Hexacosane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.16	3.98 ng	466283	Perylene-d12	15.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexacosane	366	C26H54	000630-01-3	91
2		Docosane, 11-butyl-	366	C26H54	013475-76-8	91
3		Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	91
4		Docosane	310	C22H46	000629-97-0	91
5		Heneicosane	296	C21H44	000629-94-7	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101618\
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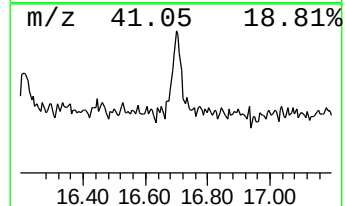
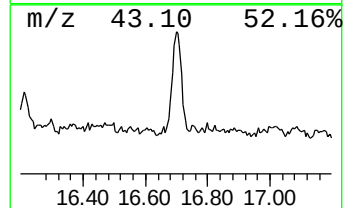
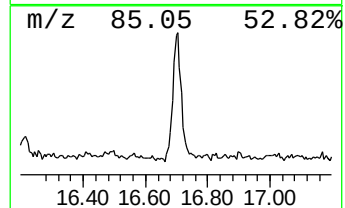
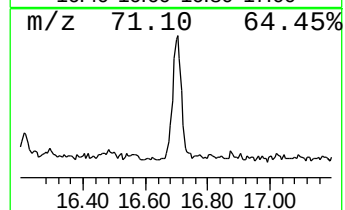
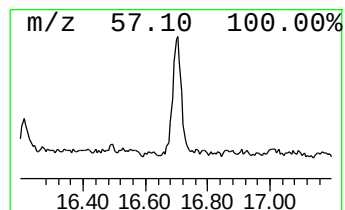
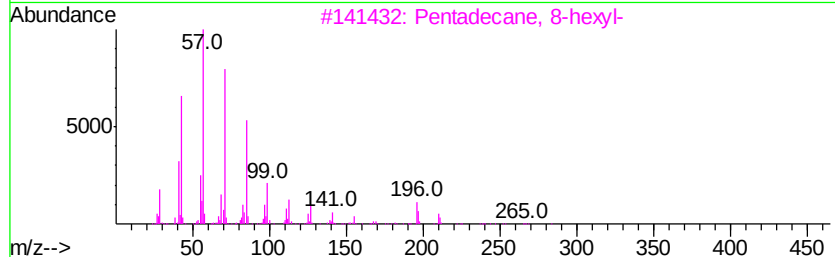
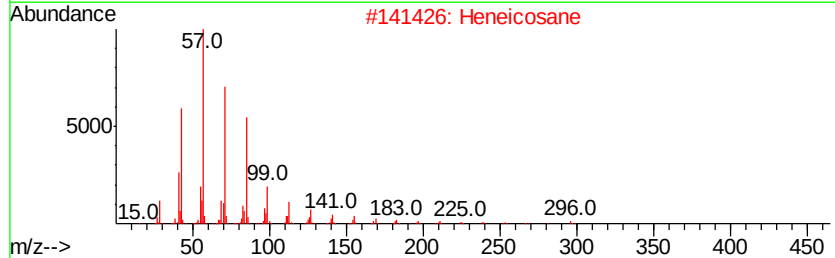
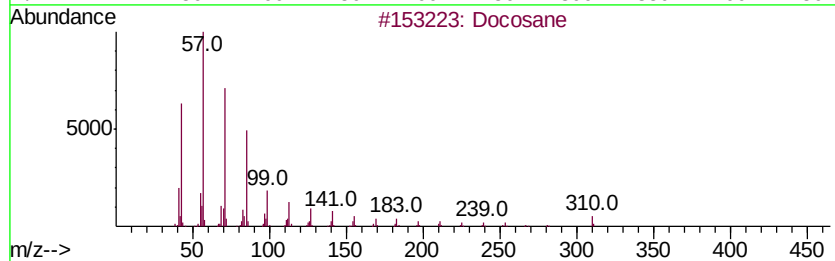
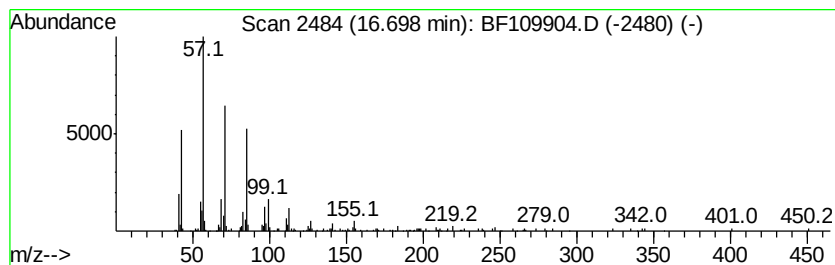
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Docosane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.70	2.85 ng	333485	Perylene-d12	15.69
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Docosane		310 C22H46	000629-97-0 90
2	Heneicosane		296 C21H44	000629-94-7 90
3	Pentadecane, 8-hexyl-		296 C21H44	013475-75-7 87
4	5,5,7,7-Tetraethylundecane		268 C19H40	1000360-42-0 86
5	Heptadecane		240 C17H36	000629-78-7 86



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF101618\
Data File : BF109904.D
Acq On : 16 Oct 2018 14:46
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Sample : J5526-10
Misc :
ALS Vial : 6 Sample Multiplier: 1

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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
Butane, 2-methoxy...	2.34	16.8	ng	1386160	1	6.98	1650360	20.0
unknown3.60	3.60	3.0	ng	245871	1	6.98	1650360	20.0
2-Pentanone, 4-hy...	5.23	4.2	ng	348538	1	6.98	1650360	20.0
unknown6.75	6.75	86.6	ng	7148990	1	6.98	1650360	20.0
1-Octadecene	12.54	2.0	ng	149787	4	11.52	1458980	20.0
Octadecanoic acid	12.80	2.2	ng	161195	4	11.52	1458980	20.0
1-Heneicosanol	13.97	6.5	ng	577675	5	14.16	1767360	20.0
Heptadecane, 2,6,...	14.61	3.5	ng	310355	5	14.16	1767360	20.0
Eicosane	14.93	3.2	ng	377214	6	15.69	2342270	20.0
Tetracosane	15.29	3.9	ng	460604	6	15.69	2342270	20.0
Hexacosane	16.16	4.0	ng	466283	6	15.69	2342270	20.0
Docosane	16.70	2.9	ng	333485	6	15.69	2342270	20.0