

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF030518\
 Data File : BF103439.D
 Acq On : 6 Mar 2018 4:50
 Operator : SJ/JU
 Sample : J1761-01
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TR-04-0302018

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:\HPCHEM1\BNA_F\METHODS\8270-BF022218.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	2.116	3	5	12	rVB	156375	182352	5.97%	0.681%
2	5.104	510	513	526	rBV	47250	71676	2.35%	0.268%
3	5.504	577	581	608	rBV	1889881	2893357	94.78%	10.811%
4	6.516	749	753	771	rBV	2206262	3052611	100.00%	11.407%
5	6.645	771	775	786	rVV	2486889	2718501	89.05%	10.158%
6	6.875	810	814	830	rVB	708467	813920	26.66%	3.041%
7	7.028	836	840	856	rBV	2108951	2084677	68.29%	7.790%
8	7.439	906	910	925	rBV	1506765	1667767	54.63%	6.232%
9	8.157	1028	1032	1050	rVB	905860	993675	32.55%	3.713%
10	9.092	1188	1191	1197	rBV	167649	142682	4.67%	0.533%
11	9.228	1211	1214	1224	rVB	2746815	2560539	83.88%	9.568%
12	9.333	1229	1232	1237	rBV	46542	43093	1.41%	0.161%
13	9.392	1239	1242	1250	rVB	49899	71723	2.35%	0.268%
14	9.622	1277	1281	1284	rBV2	124316	140750	4.61%	0.526%
15	9.645	1284	1285	1292	rVB4	33783	35701	1.17%	0.133%
16	9.704	1292	1295	1300	rBB3	31751	41967	1.37%	0.157%
17	9.916	1326	1331	1342	rVB	991672	989874	32.43%	3.699%
18	10.298	1394	1396	1398	rBV	68200	57569	1.89%	0.215%
19	10.322	1398	1400	1404	rVB2	58748	53526	1.75%	0.200%
20	10.545	1433	1438	1441	rBV2	37763	48957	1.60%	0.183%
21	10.710	1462	1466	1475	rBV	1182953	1269106	41.57%	4.742%
22	10.810	1479	1483	1487	rVV	63763	78019	2.56%	0.292%
23	10.904	1496	1499	1503	rVB	38405	35056	1.15%	0.131%
24	11.251	1555	1558	1560	rBV	32041	33643	1.10%	0.126%
25	11.280	1560	1563	1570	rVB2	47937	58464	1.92%	0.218%
26	11.339	1570	1573	1578	rBV3	57906	61934	2.03%	0.231%
27	11.410	1581	1585	1589	rBV	836981	814887	26.69%	3.045%
28	11.557	1607	1610	1613	rBV	193788	161975	5.31%	0.605%
29	11.616	1615	1620	1626	rBV4	31132	71449	2.34%	0.267%
30	11.674	1627	1630	1632	rBV	36991	31436	1.03%	0.117%
31	11.780	1644	1648	1651	rVB2	281819	238303	7.81%	0.890%
32	11.916	1668	1671	1674	rBV3	53188	58056	1.90%	0.217%
33	11.957	1674	1678	1681	rVB	614145	547394	17.93%	2.045%
34	12.027	1687	1690	1693	rBV2	31590	32109	1.05%	0.120%

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Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF022218.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	12.192	1715	1718	1721	rBV	51845	46498	1.52%	0.174%
36	12.469	1761	1765	1766	rBV	632954	542634	17.78%	2.028%
37	12.486	1766	1768	1776	rVB	612751	573926	18.80%	2.145%
38	12.574	1778	1783	1788	rBV	112345	114452	3.75%	0.428%
39	12.633	1790	1793	1796	rVB	27832	30608	1.00%	0.114%
40	12.792	1818	1820	1826	rVB3	28319	36888	1.21%	0.138%
41	12.845	1826	1829	1831	rBV	107702	91965	3.01%	0.344%
42	12.992	1850	1854	1858	rBV	1746984	1719691	56.34%	6.426%
43	13.204	1887	1890	1893	rVB	139321	120905	3.96%	0.452%
44	13.427	1925	1928	1930	rBV3	29551	33250	1.09%	0.124%
45	13.545	1945	1948	1953	rVB	126164	108292	3.55%	0.405%
46	13.874	2001	2004	2012	rBV	197123	221641	7.26%	0.828%
47	14.068	2033	2037	2045	rVB	523060	567472	18.59%	2.120%
48	14.198	2056	2059	2062	rBV	59840	65656	2.15%	0.245%
49	15.580	2291	2294	2301	rVB2	259228	331398	10.86%	1.238%

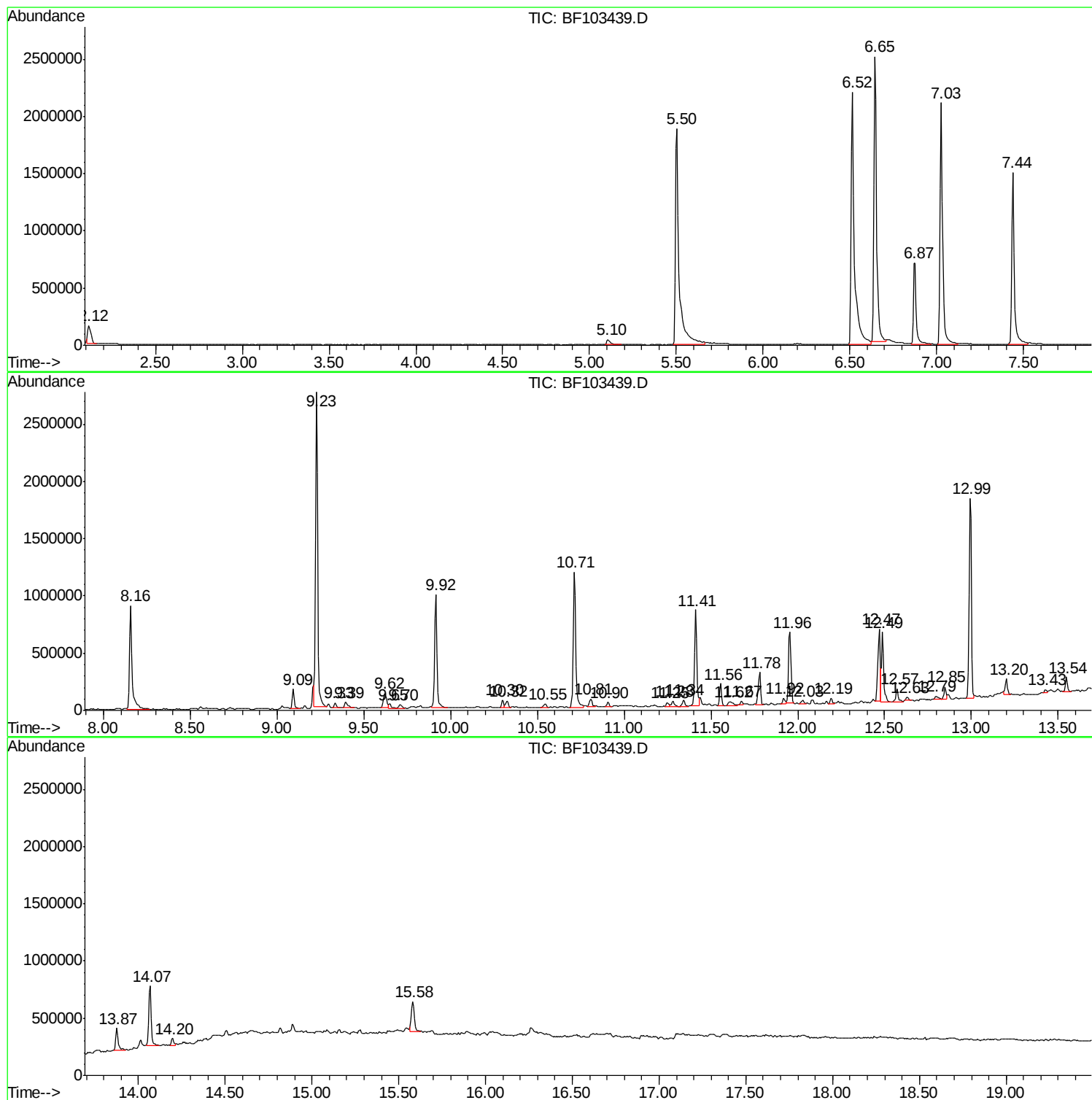
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF022218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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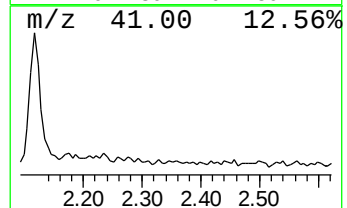
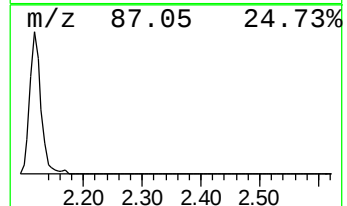
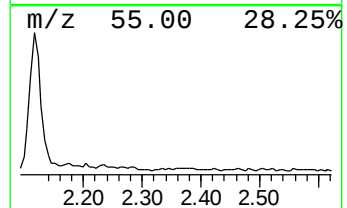
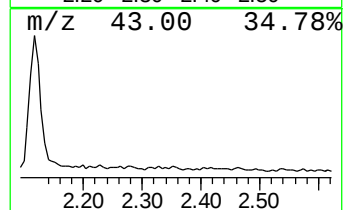
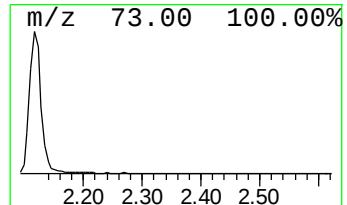
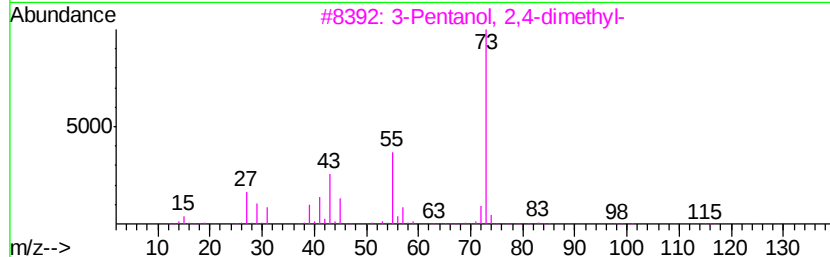
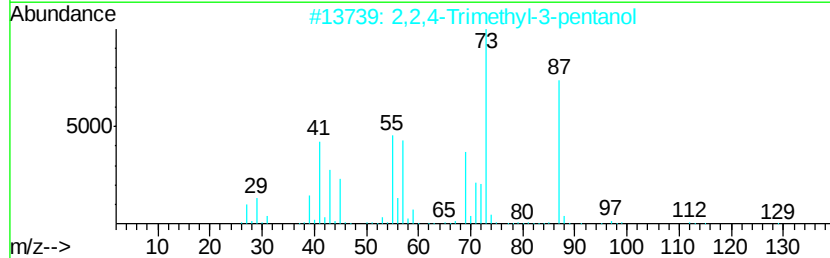
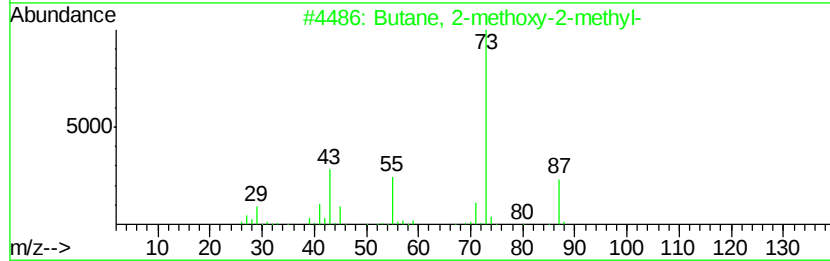
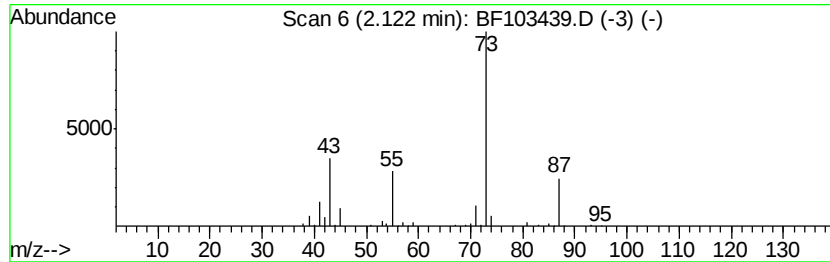
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.12	4.48 ng	182352	1,4-Dichlorobenzene-d4	6.87

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
3			3-Pentanol, 2,4-dimethyl-	116	C7H16O	000600-36-2	37
4			Silane, ethyltrimethyl-	102	C5H14Si	003439-38-1	33
5			Silane, tetramethyl-	88	C4H12Si	000075-76-3	16



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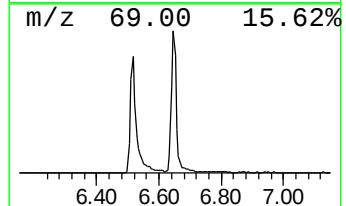
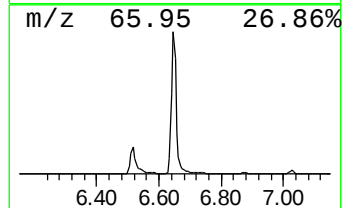
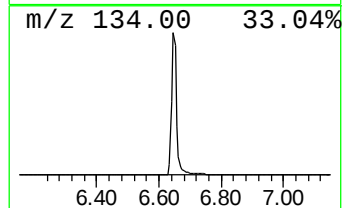
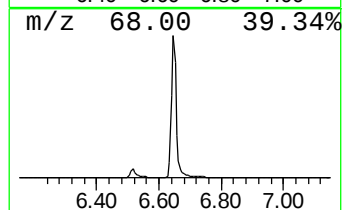
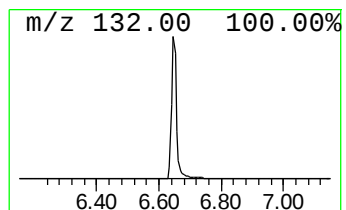
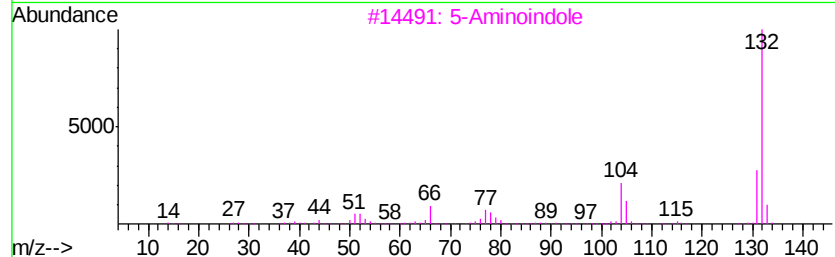
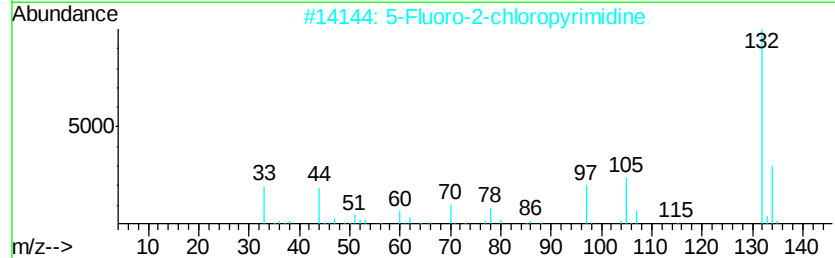
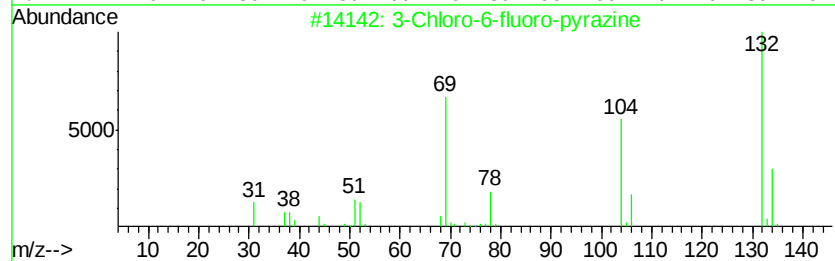
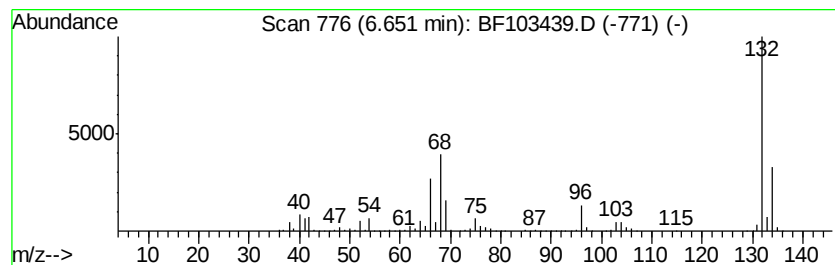
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown6.65 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.65	66.80 ng	2718500	1,4-Dichlorobenzene-d4	6.87

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	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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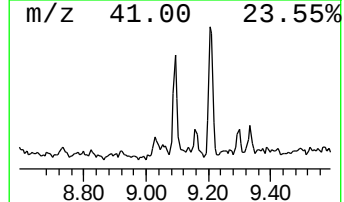
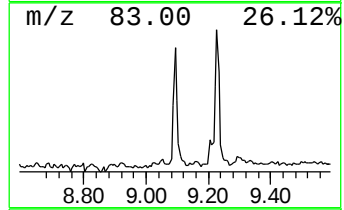
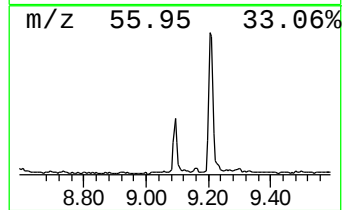
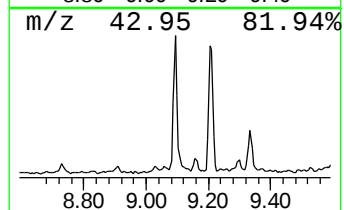
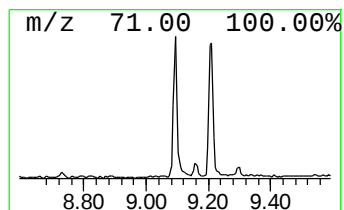
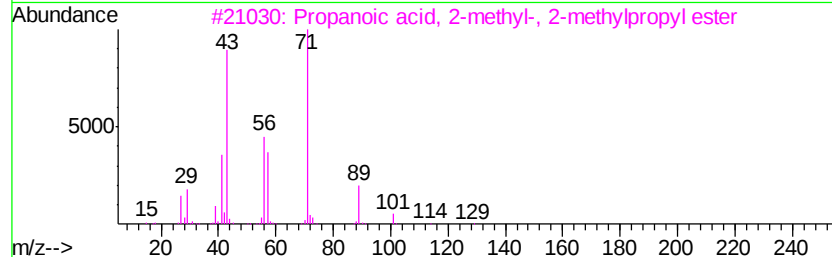
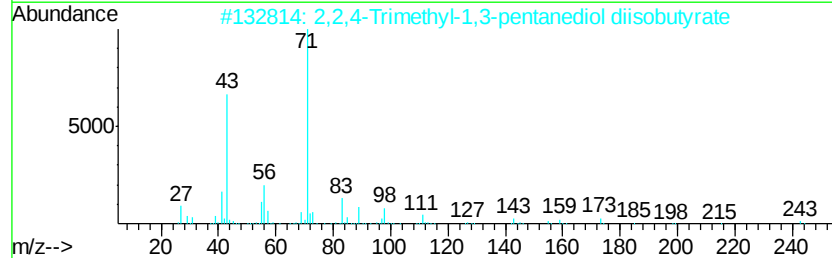
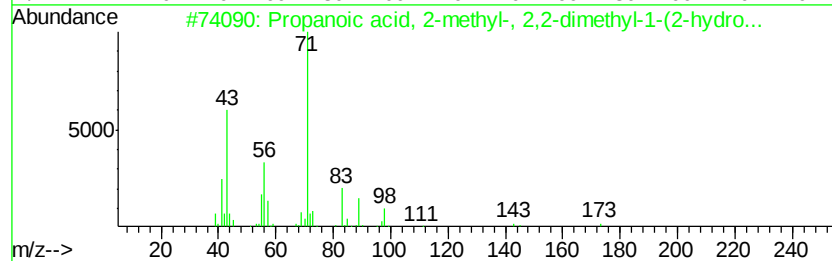
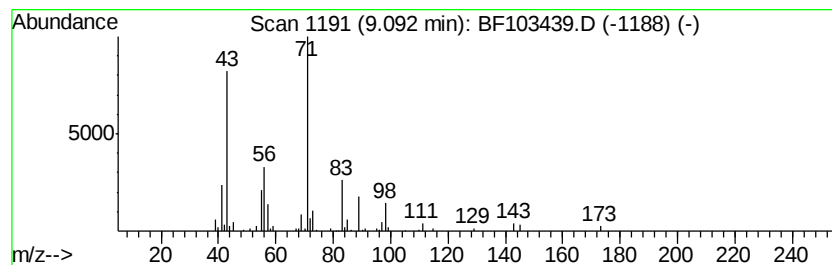
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Propanoic acid, 2-methyl-, ... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.09	2.88 ng	142682	Acenaphthene-d10	9.92

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propanoic acid, 2-methyl-, 2,2-d...	216	C12H24O3	074367-33-2	90
2		2,2,4-Trimethyl-1,3-pentanediol ...	286	C16H30O4	006846-50-0	64
3		Propanoic acid, 2-methyl-, 2-met...	144	C8H16O2	000097-85-8	50
4		1-Hexadecen-3-ol, 3,5,11,15-tetr...	296	C20H40O	1000262-55-5	43
5		Isophytol	296	C20H40O	000505-32-8	43



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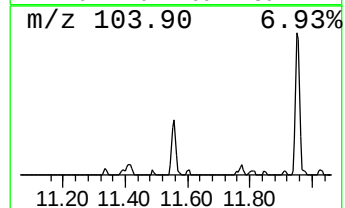
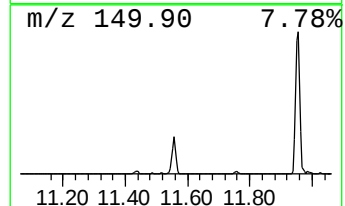
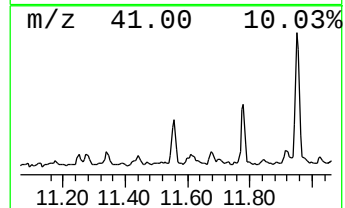
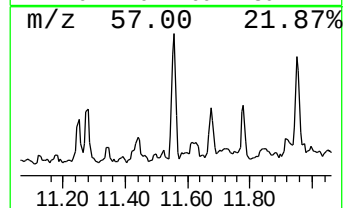
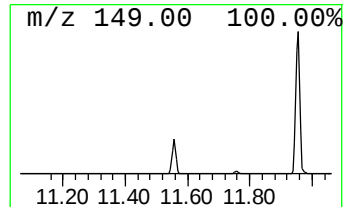
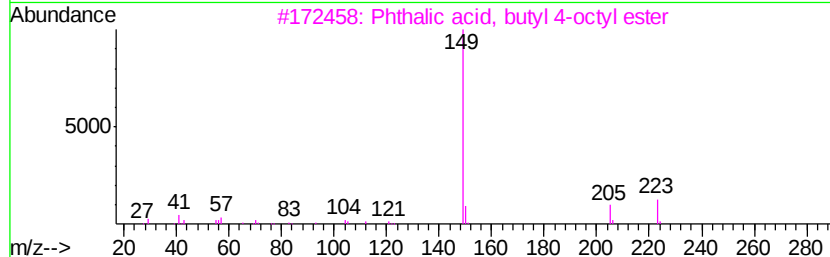
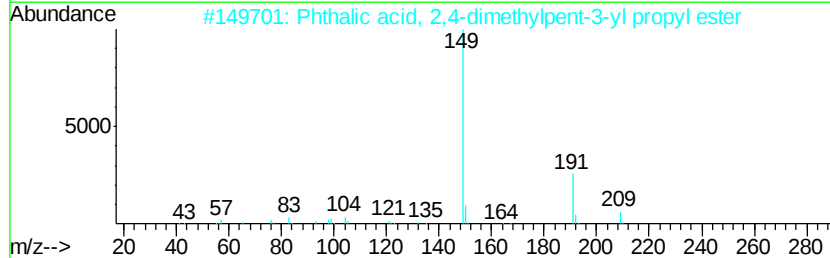
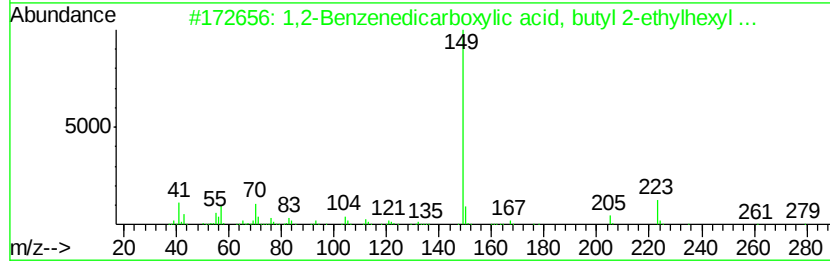
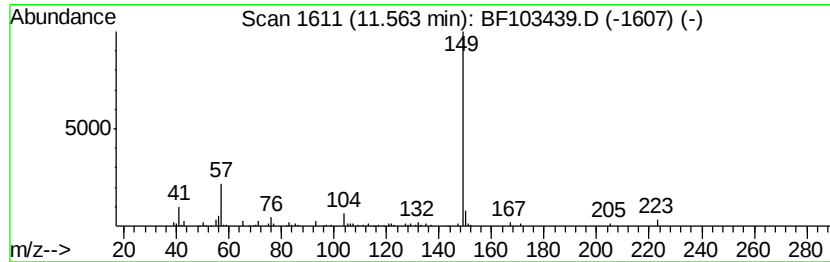
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 5 1,2-Benzenedicarboxylic aci... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.56	3.98 ng	161975	Phenanthrene-d10	11.41

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,2-Benzenedicarboxylic acid, bu...	334	C20H30O4	000085-69-8	83	
2	Phthalic acid, 2,4-dimethylpent-...	306	C18H26O4	1000356-84-2	78	
3	Phthalic acid, butyl 4-octyl ester	334	C20H30O4	1000314-83-8	78	
4	1,2-Benzenedicarboxylic acid, bi...	278	C16H22O4	000084-69-5	78	
5	Phthalic acid, isobutyl propyl e...	264	C15H20O4	1000308-97-1	78	



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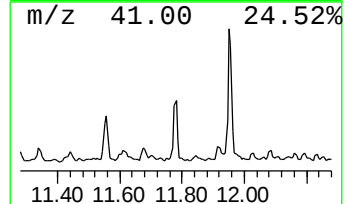
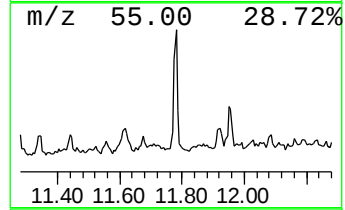
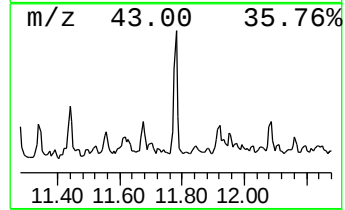
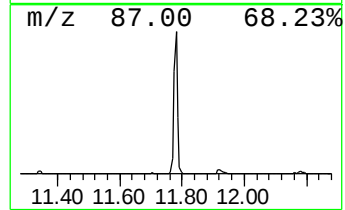
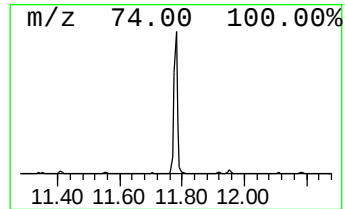
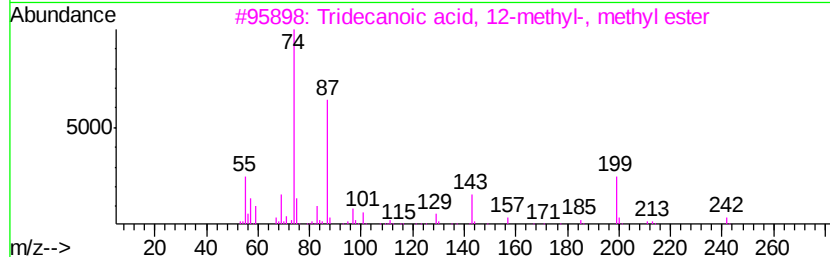
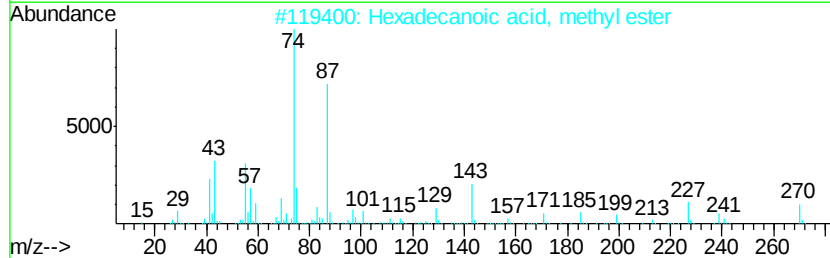
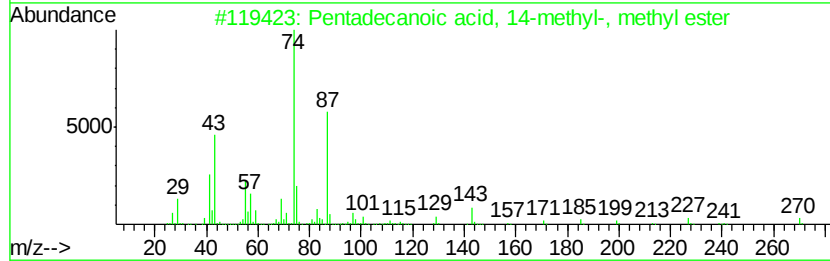
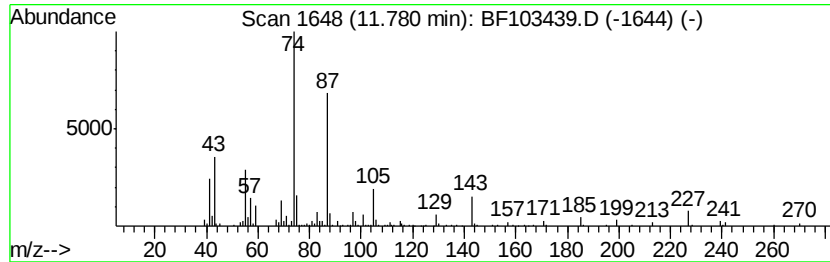
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Pentadecanoic acid, 14-meth... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.78	5.85 ng	238303	Phenanthrene-d10	11.41	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	83
2	Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	83
3	Tridecanoic acid, 12-methyl-, me...	242	C15H30O2	005129-58-8	83
4	Pentadecanoic acid, methyl ester	256	C16H32O2	007132-64-1	80
5	Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	80



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Data File : BF103439.D
Acq On : 6 Mar 2018 4:50
Operator : SJ/JU
Sample : J1761-01
Misc :
ALS Vial : 27 Sample Multiplier: 1

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ClientSampleId :
TR-04-0302018

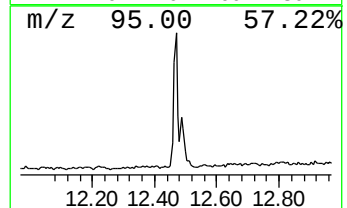
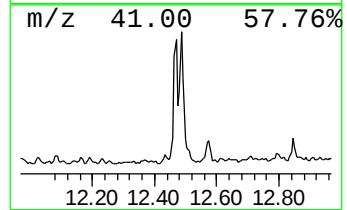
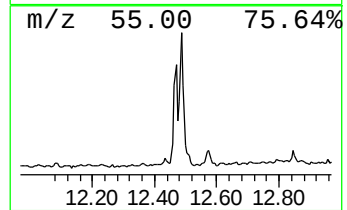
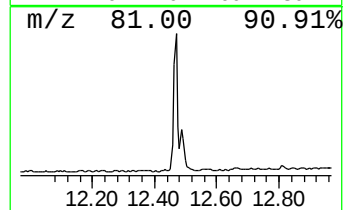
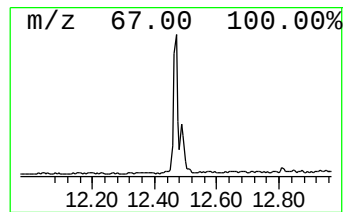
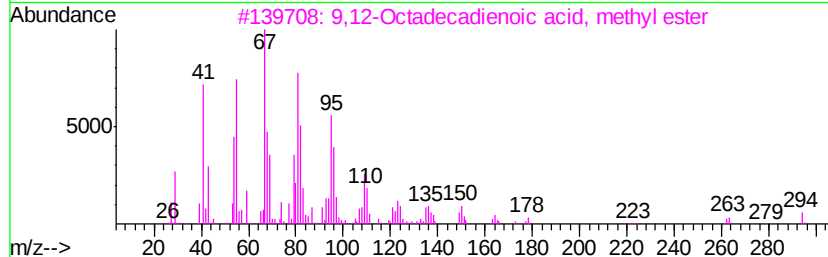
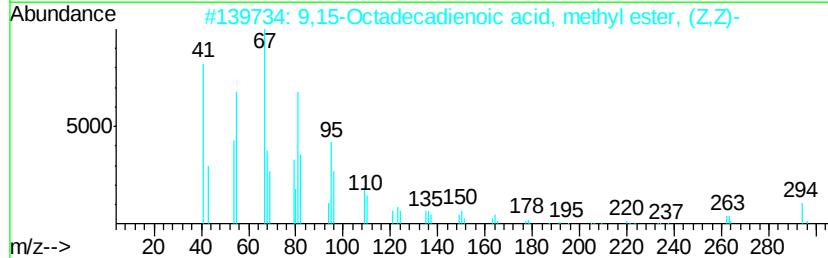
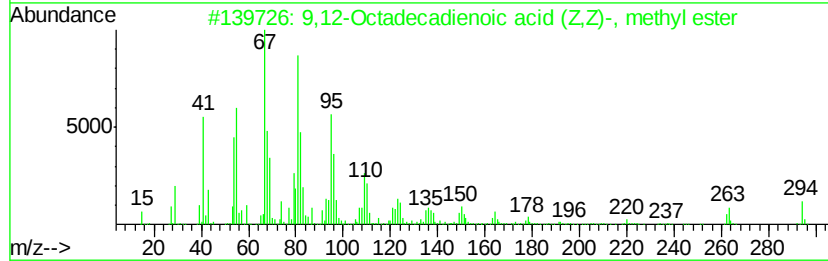
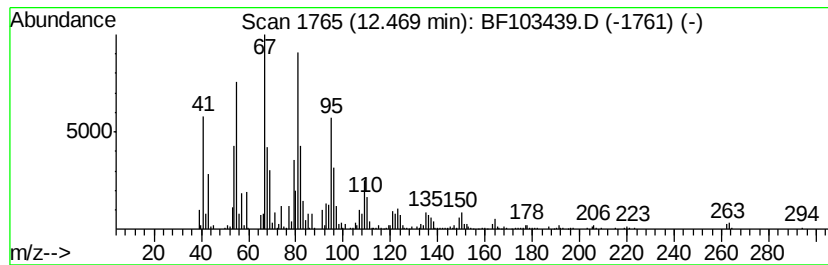
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 9,12-Octadecadienoic acid (... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.47	13.32 ng	542634	Phenanthrene-d10	11.41

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9,12-Octadecadienoic acid (Z,Z)-...	294	C19H34O2	000112-63-0	99
2		9,15-Octadecadienoic acid, methy...	294	C19H34O2	017309-05-6	99
3		9,12-Octadecadienoic acid, methv...	294	C19H34O2	002462-85-3	99
4		9,12-Octadecadienoic acid (Z,Z)-	280	C18H32O2	000060-33-3	95
5		2-Chloroethyl linoleate	342	C20H35ClO2	025525-76-2	94



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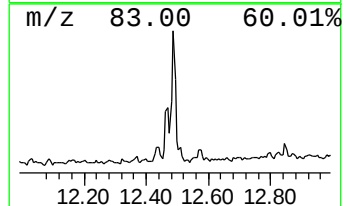
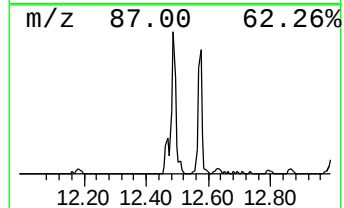
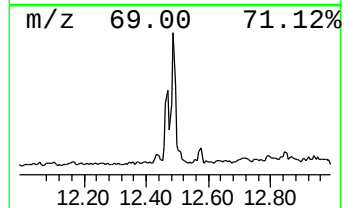
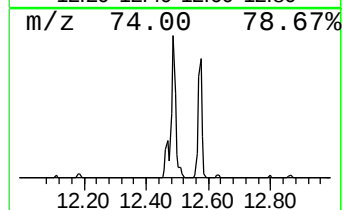
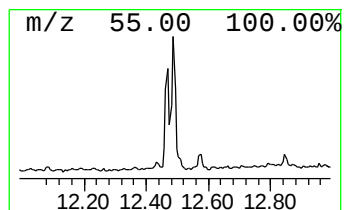
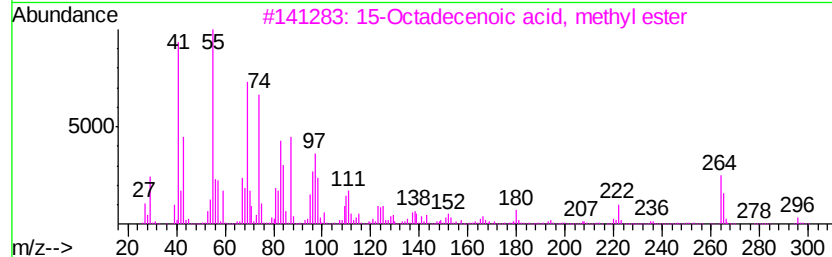
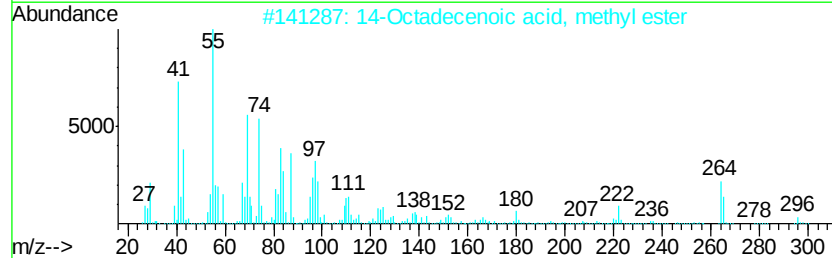
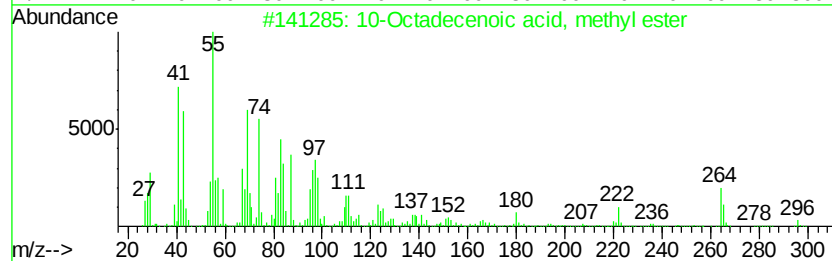
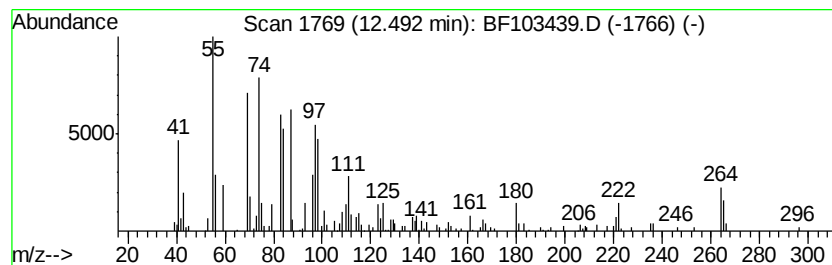
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 10-Octadecenoic acid, methyl... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.49	14.09 ng	573926	Phenanthrene-d10	11.41

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		10-Octadecenoic acid, methyl ester	296	C19H36O2	013481-95-3	91
2		14-Octadecenoic acid, methyl ester	296	C19H36O2	056554-48-4	91
3		15-Octadecenoic acid, methyl ester	296	C19H36O2	004764-72-1	91
4		9-Octadecenoic acid (Z)-, methyl...	296	C19H36O2	000112-62-9	87
5		16-Octadecenoic acid, methyl ester	296	C19H36O2	056554-49-5	86



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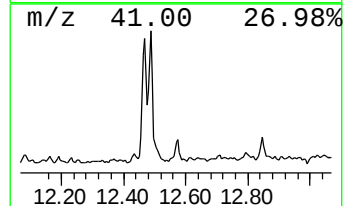
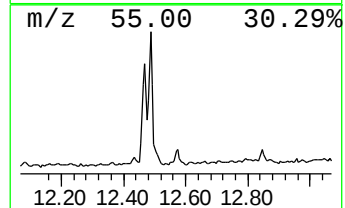
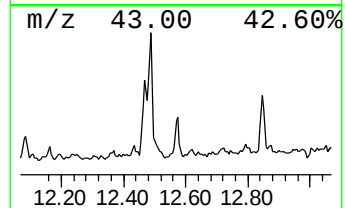
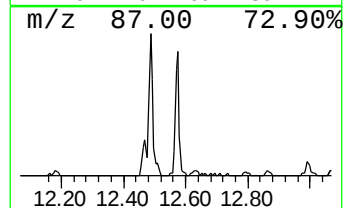
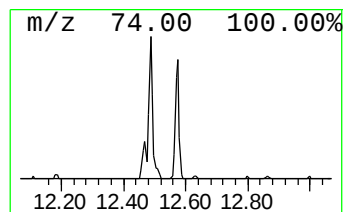
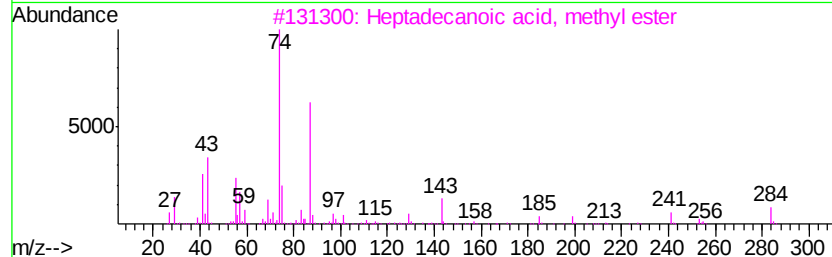
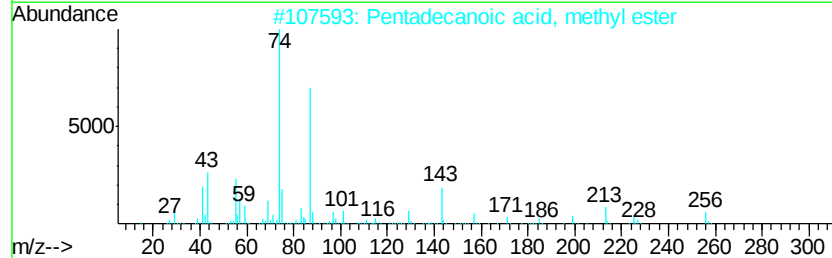
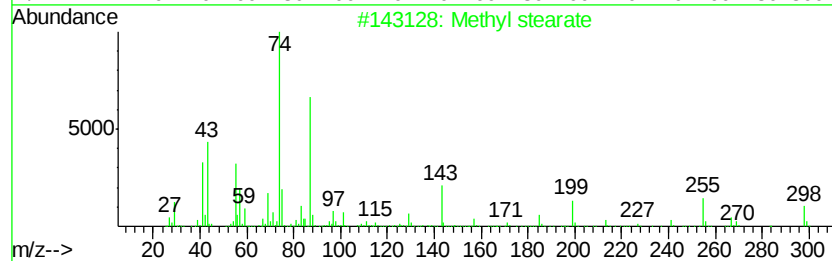
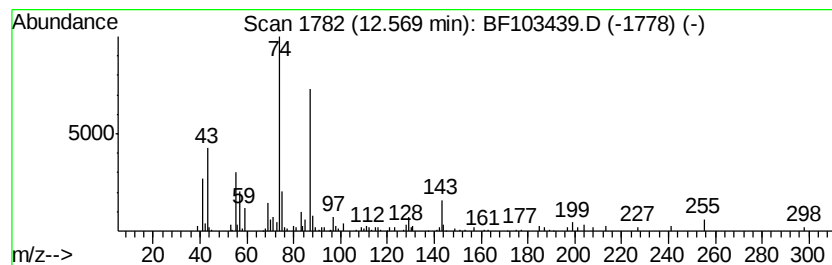
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Methyl stearate Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.57	2.81 ng	114452	Phenanthrene-d10	11.41

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Methyl stearate	298	C19H38O2	000112-61-8	94
2		Pentadecanoic acid, methyl ester	256	C16H32O2	007132-64-1	91
3		Heptadecanoic acid, methyl ester	284	C18H36O2	001731-92-6	91
4		Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	90
5		Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	90



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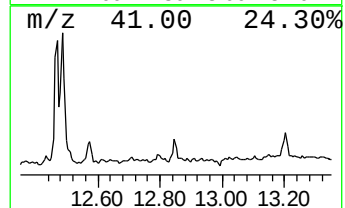
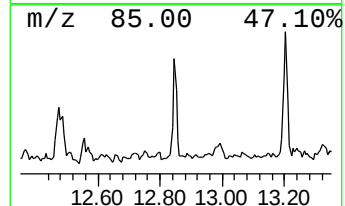
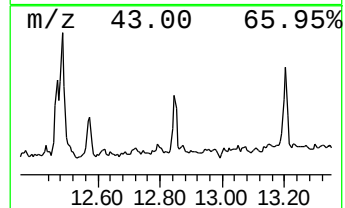
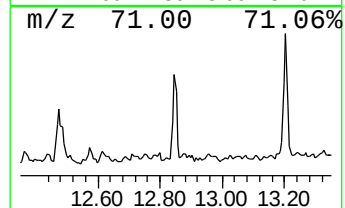
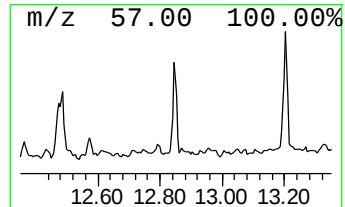
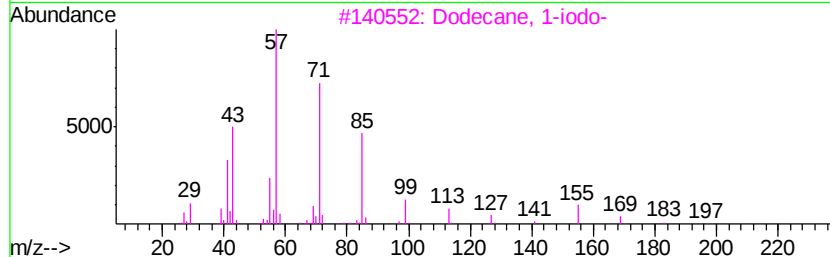
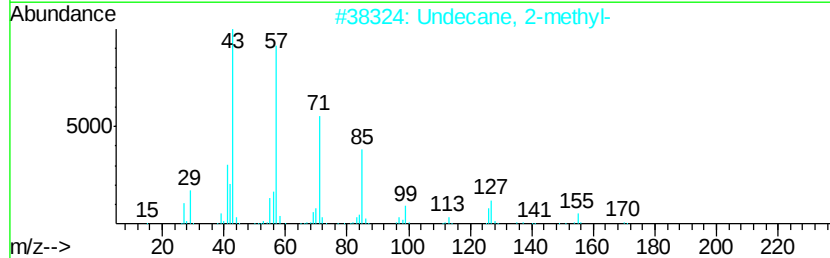
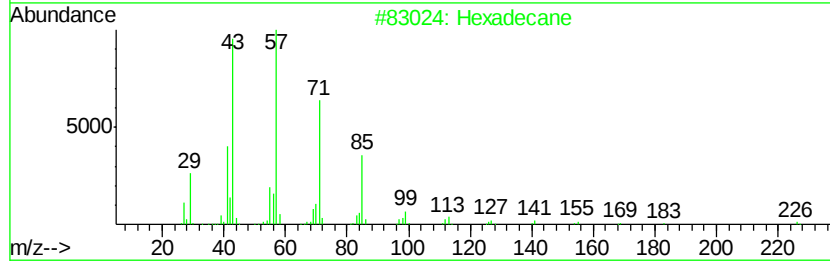
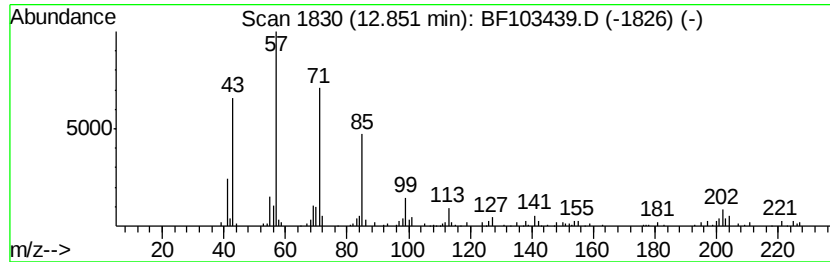
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Hexadecane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD		R.T.
12.85	3.24 ng	91965	Chrysene-d12		14.07
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	87
2	Undecane, 2-methyl-	170	C12H26	007045-71-8	80
3	Dodecane, 1-iodo-	296	C12H25I	004292-19-7	80
4	Dodecane, 2-methyl-	184	C13H28	001560-97-0	80
5	Heptadecane	240	C17H36	000629-78-7	80



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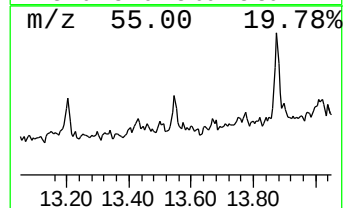
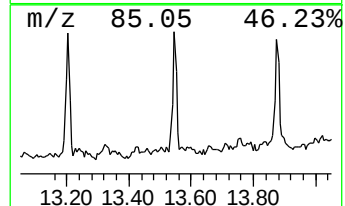
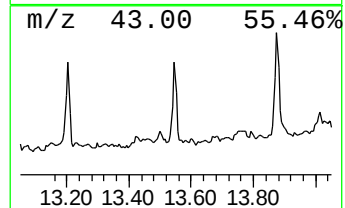
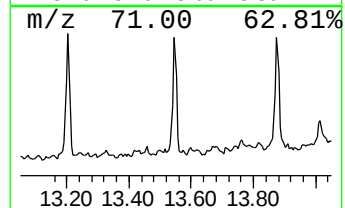
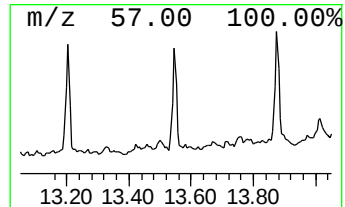
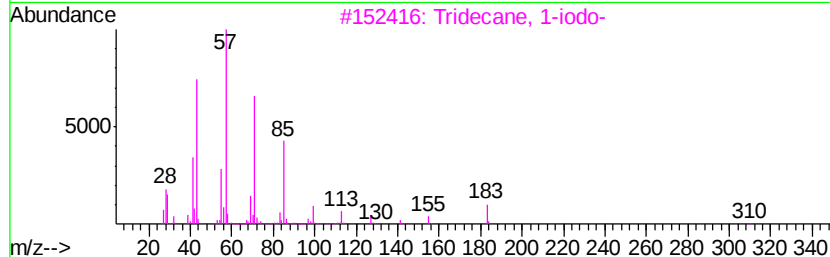
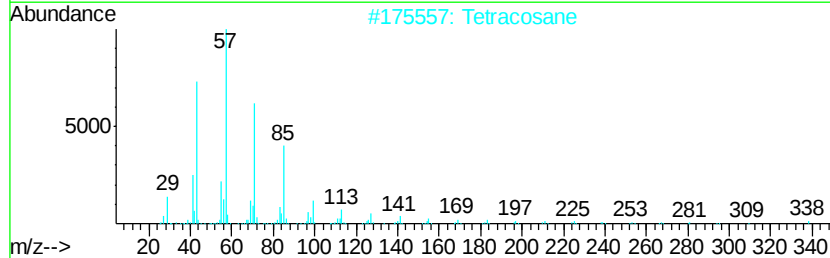
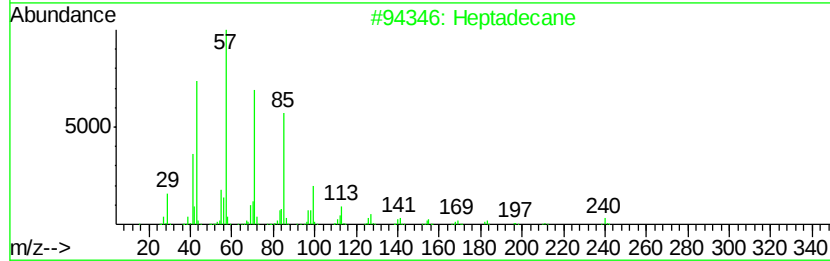
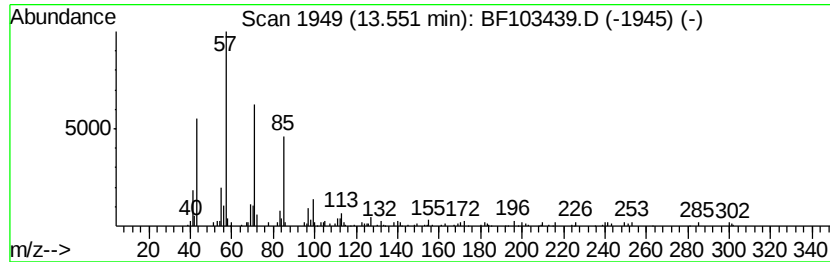
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Heptadecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.55	3.82 ng	108292	Chrysene-d12	14.07	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane	240	C17H36	000629-78-7	95
2	Tetracosane	338	C24H50	000646-31-1	83
3	Tridecane, 1-iodo-	310	C13H27I	035599-77-0	83
4	Heptacosane	380	C27H56	000593-49-7	83
5	Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	80



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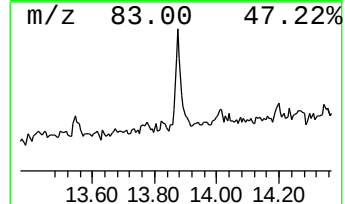
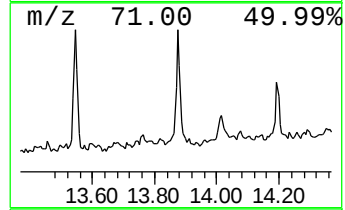
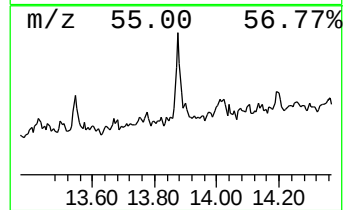
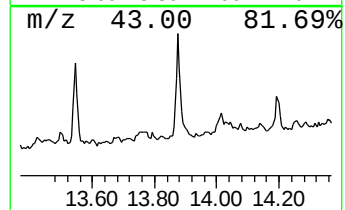
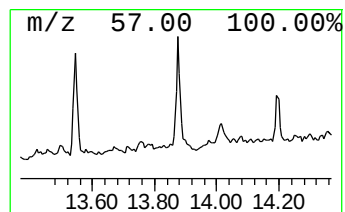
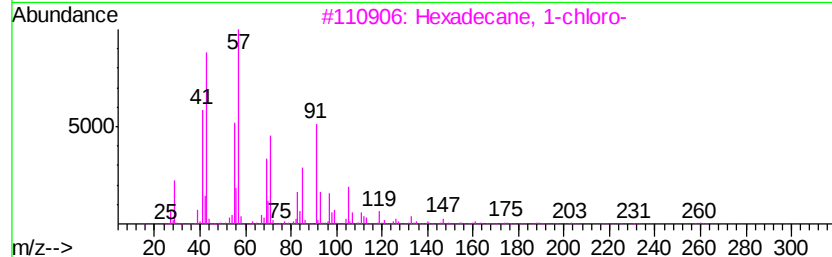
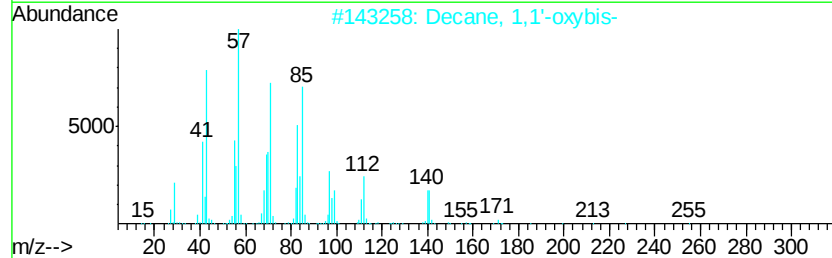
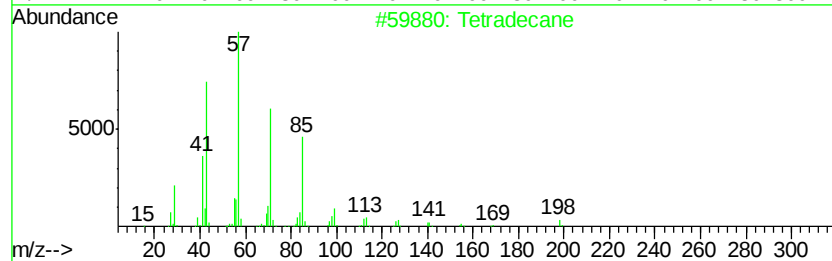
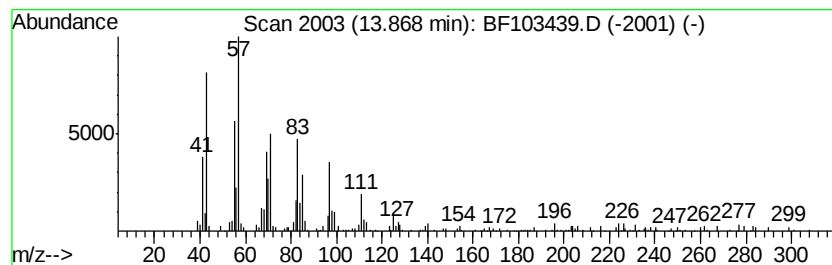
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 Tetradecane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.87	7.81 ng	221641	Chrysene-d12	14.07

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetradecane	198	C14H30	000629-59-4	90
2			Decane, 1,1'-oxybis-	298	C20H42O	002456-28-2	83
3			Hexadecane, 1-chloro-	260	C16H33Cl	004860-03-1	72
4			Oxalic acid, isobutyl tetradecyl...	342	C20H38O4	1000309-37-9	68
5			Oxalic acid, isobutyl pentadecyl...	356	C21H40O4	1000309-38-0	68



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF030518\
Data File : BF103439.D
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Sample : J1761-01
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
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ClientSampleId :
TR-04-0302018

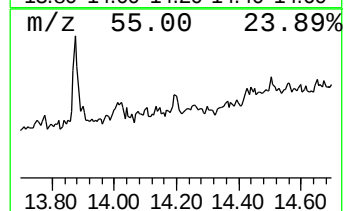
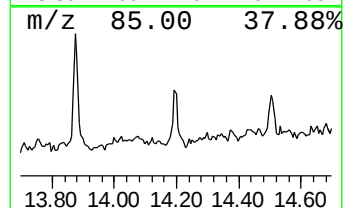
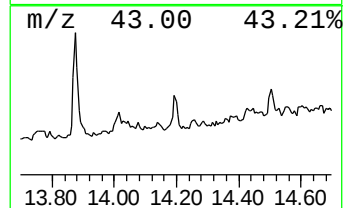
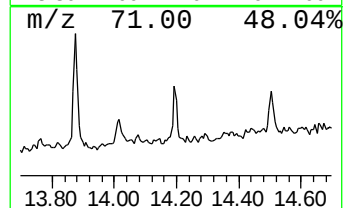
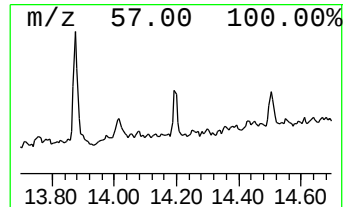
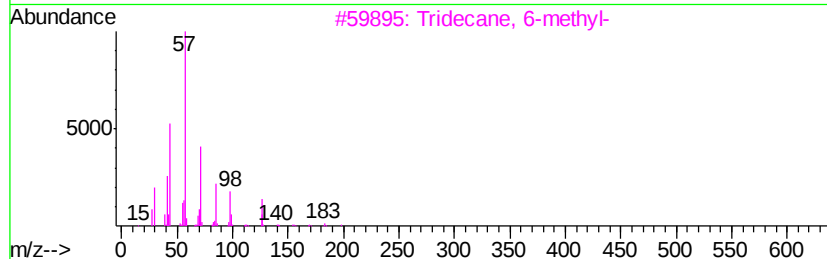
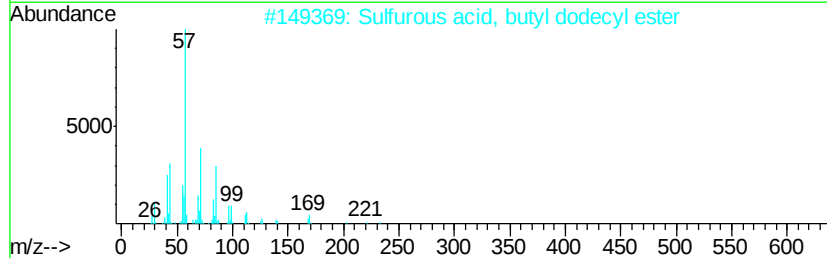
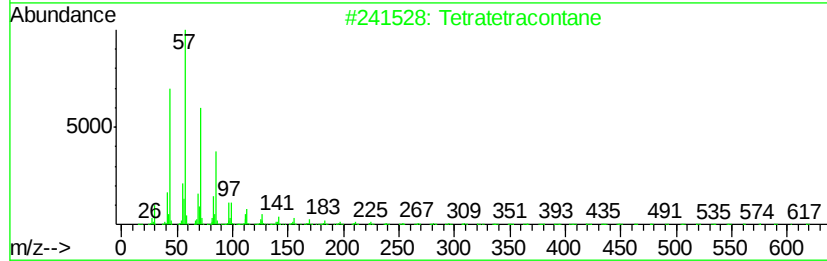
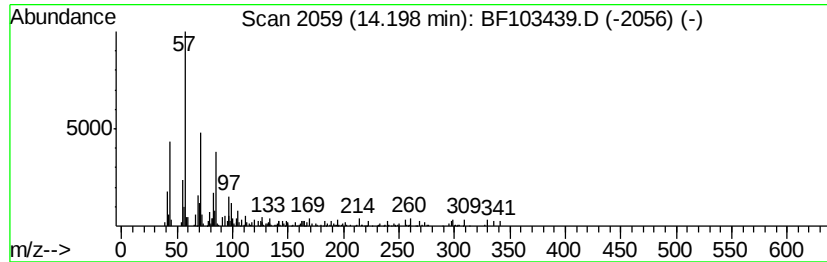
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Tetratetracontane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.20	2.31 ng	65656	Chrysene-d12	14.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetratetracontane	619	C44H90	007098-22-8	74
2		Sulfurous acid, butyl dodecyl ester	306	C16H34O3S	1000309-17-9	64
3		Tridecane, 6-methyl-	198	C14H30	013287-21-3	64
4		Heptacosane, 1-chloro-	414	C27H55Cl	062016-79-9	64
5		Ethanol, 2-(dodecyloxy)-	230	C14H30O2	004536-30-5	64



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF030518\
 Data File : BF103439.D
 Acq On : 6 Mar 2018 4:50
 Operator : SJ/JU
 Sample : J1761-01
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TR-04-0302018

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF022218.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
Butane, 2-methoxy...	2.12	4.5	ng	182352	1	6.87	813920	20.0
unknown6.65	6.65	66.8	ng	2718500	1	6.87	813920	20.0
Propanoic acid, 2...	9.09	2.9	ng	142682	3	9.92	989874	20.0
1,2-Benzenedicarb...	11.56	4.0	ng	161975	4	11.41	814887	20.0
Pentadecanoic aci...	11.78	5.8	ng	238303	4	11.41	814887	20.0
9,12-Octadecadien...	12.47	13.3	ng	542634	4	11.41	814887	20.0
10-Octadecenoic a...	12.49	14.1	ng	573926	4	11.41	814887	20.0
Methyl stearate	12.57	2.8	ng	114452	4	11.41	814887	20.0
Hexadecane	12.85	3.2	ng	91965	5	14.07	567472	20.0
Heptadecane	13.55	3.8	ng	108292	5	14.07	567472	20.0
Tetradecane	13.87	7.8	ng	221641	5	14.07	567472	20.0
Tetratetracontane	14.20	2.3	ng	65656	5	14.07	567472	20.0