

Data Path : Z:\HPCHEM1\BNA F\DATA\BF032318\
 Data File : BF103884.D
 Acq On : 23 Mar 2018 17:00
 Operator : JU/SJ
 Sample : J1748-09
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 AU-05-032218-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:\HPCHEM1\BNA_F\METHODS\8270-BF031518.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	2.304	31	37	47	rVB	109775	146558	3.13%	0.396%
2	5.087	508	510	521	rBV	47143	67212	1.44%	0.182%
3	5.216	527	532	541	rVB	134212	147216	3.14%	0.398%
4	5.604	593	598	601	rBV	2938124	2879350	61.48%	7.784%
5	6.598	762	767	779	rBV	2670376	3041936	64.95%	8.223%
6	6.745	788	792	795	rBV	3263086	3019713	64.47%	8.163%
7	6.975	827	831	834	rBV	1126606	873653	18.65%	2.362%
8	7.133	854	858	861	rBV	2703885	2303594	49.18%	6.227%
9	7.539	922	927	930	rBV	1930582	1770128	37.79%	4.785%
10	8.257	1045	1049	1052	rBV	1412509	1112420	23.75%	3.007%
11	9.333	1227	1232	1235	rBV	3459888	3073296	65.62%	8.308%
12	9.398	1240	1243	1246	rVV	82080	61666	1.32%	0.167%
13	9.722	1292	1298	1301	rBV	297700	257117	5.49%	0.695%
14	9.927	1330	1333	1337	rVB	150719	124889	2.67%	0.338%
15	10.016	1345	1348	1352	rBV	1321836	1145263	24.45%	3.096%
16	10.427	1416	1418	1421	rVB	149575	126127	2.69%	0.341%
17	10.651	1450	1456	1458	rBV	51325	60801	1.30%	0.164%
18	10.810	1479	1483	1486	rBV	1864629	1662580	35.50%	4.495%
19	10.904	1496	1499	1509	rVB	131796	159101	3.40%	0.430%
20	11.351	1572	1575	1578	rBV	60674	54791	1.17%	0.148%
21	11.510	1598	1602	1605	rBV	1157163	1011875	21.60%	2.735%
22	11.533	1605	1606	1610	rVB	86580	59810	1.28%	0.162%
23	11.886	1662	1666	1669	rBV	1880193	1490585	31.83%	4.030%
24	12.015	1685	1688	1691	rBV2	71092	73851	1.58%	0.200%
25	12.580	1779	1784	1786	rBV	4242537	4683621	100.00%	12.662%
26	12.598	1786	1787	1794	rVV2	3305236	2635653	56.27%	7.125%
27	12.680	1797	1801	1804	rVV	762327	612580	13.08%	1.656%
28	12.727	1806	1809	1812	rBV	88268	75864	1.62%	0.205%
29	12.915	1837	1841	1844	rBV4	44311	61049	1.30%	0.165%
30	12.957	1845	1848	1855	rVB2	86130	134895	2.88%	0.365%
31	13.098	1868	1872	1875	rBV	2486401	2200605	46.99%	5.949%
32	13.974	2019	2021	2025	rBV	377293	354693	7.57%	0.959%
33	14.168	2050	2054	2057	rBV	827128	808773	17.27%	2.186%
34	15.715	2313	2317	2323	rVB	504849	699615	14.94%	1.891%

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Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
AU-05-032218-B

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-BF031518.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

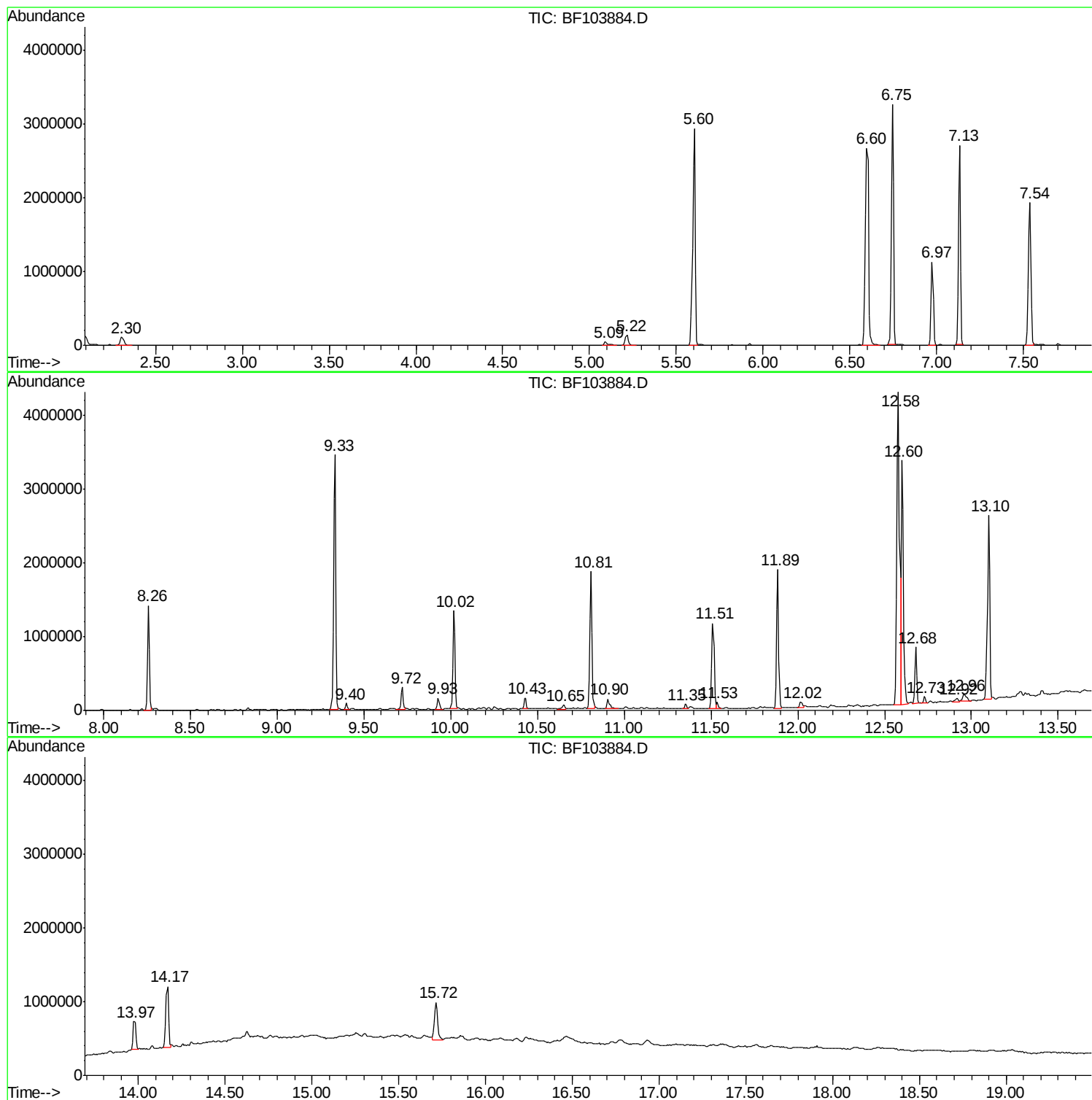
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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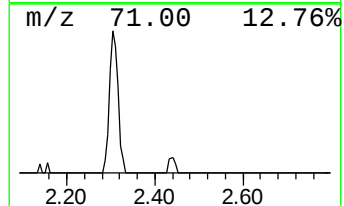
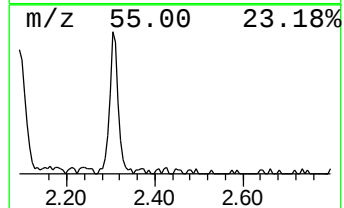
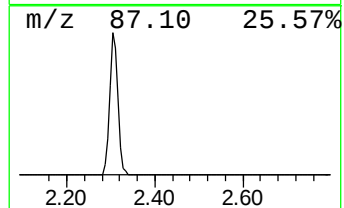
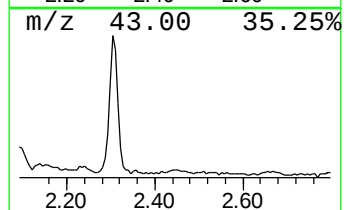
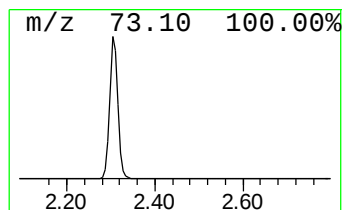
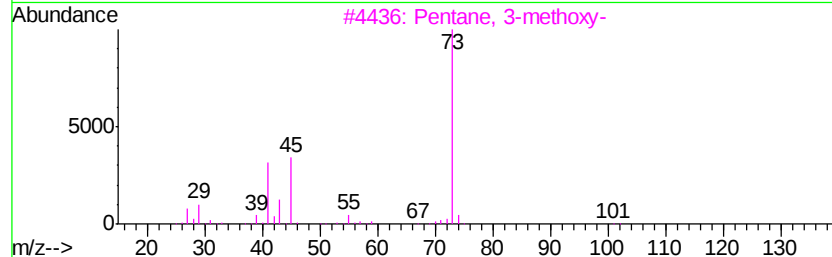
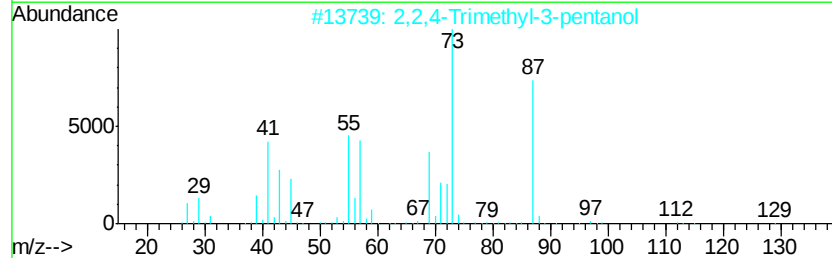
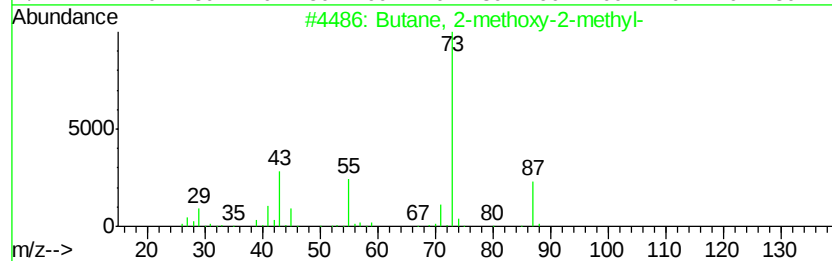
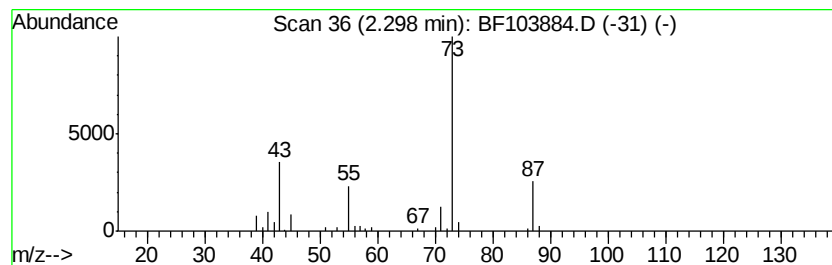
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.30	3.36 ng	146558	1,4-Dichlorobenzene-d4	6.97

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
3			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
4			3-Hexanol, 2-methyl-	116	C7H16O	000617-29-8	17
5			Boronic acid, ethyl-, dimethyl e...	102	C4H11BO2	007318-82-3	10



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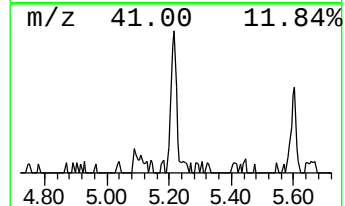
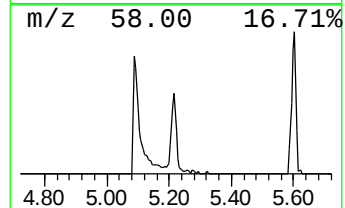
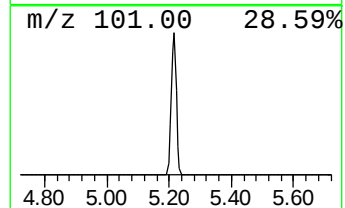
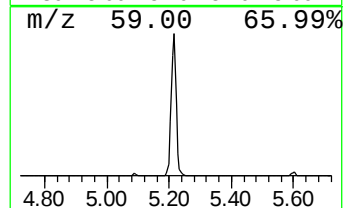
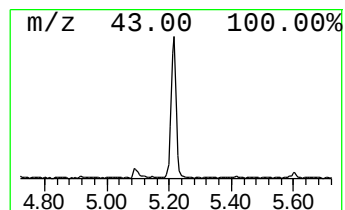
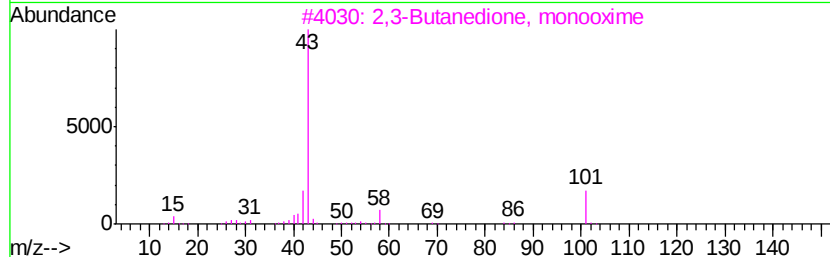
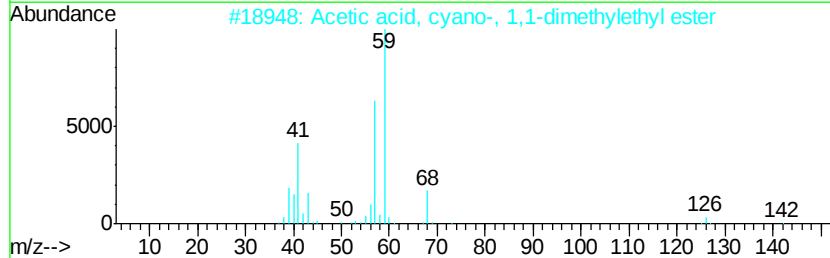
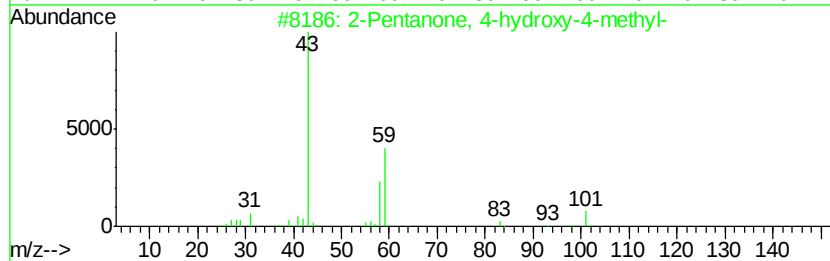
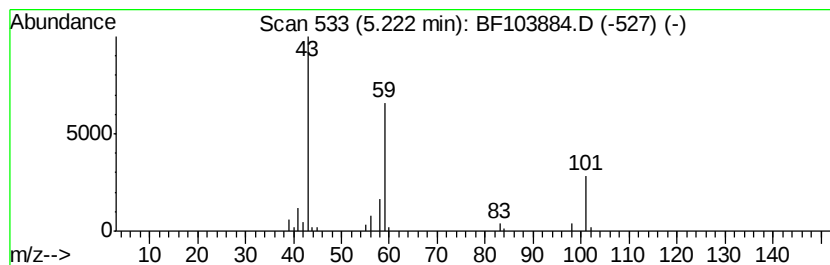
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.22	3.37 ng	147216	1,4-Dichlorobenzene-d4	6.97

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, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	17
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
4			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5			5-Hexen-2-one	98	C6H10O	000109-49-9	9



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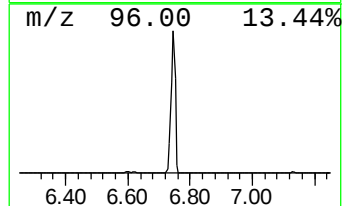
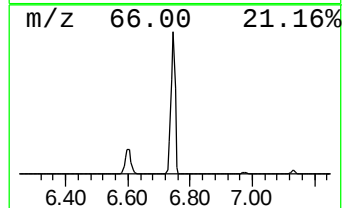
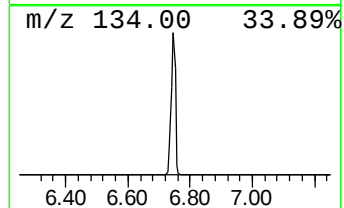
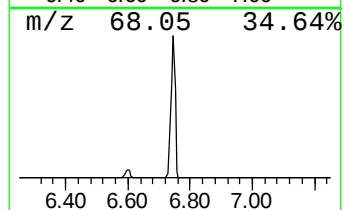
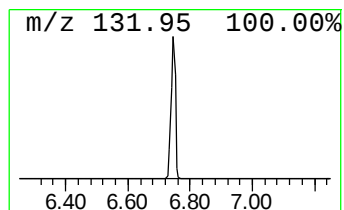
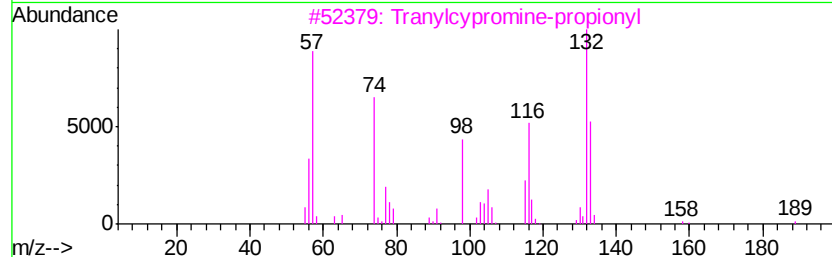
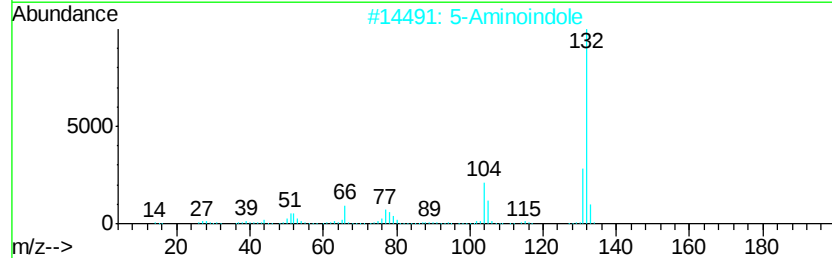
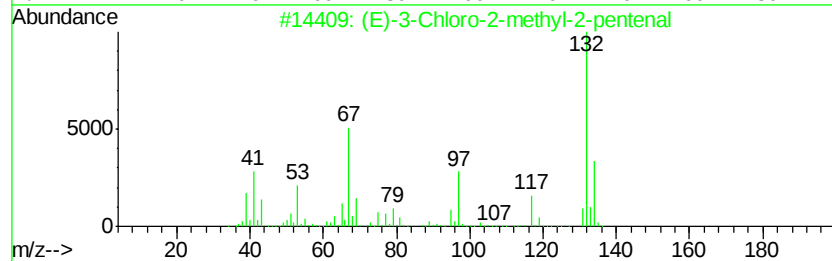
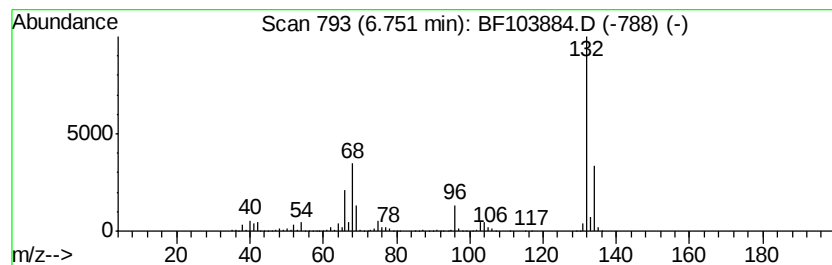
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown6.75 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.75	69.13 ng	3019710	1,4-Dichlorobenzene-d4	6.97

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
2			5-Aminoindole	132	C8H8N2	005192-03-0	12
3			Tranvlcyproline-propionyl	189	C12H15NO	1000123-86-3	10
4			4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9
5			4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	9



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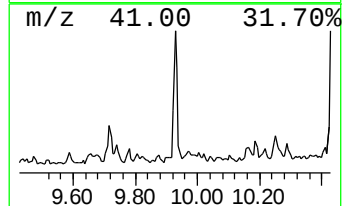
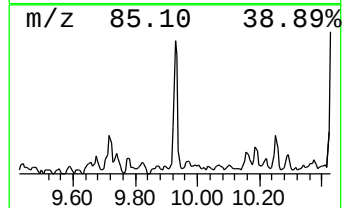
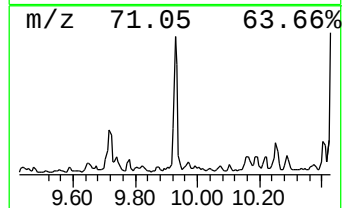
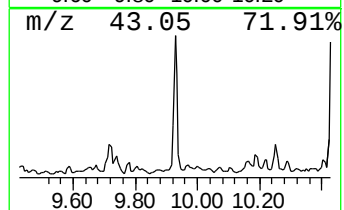
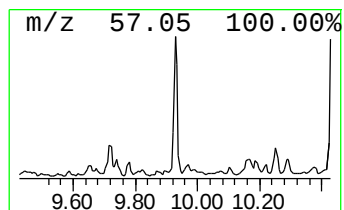
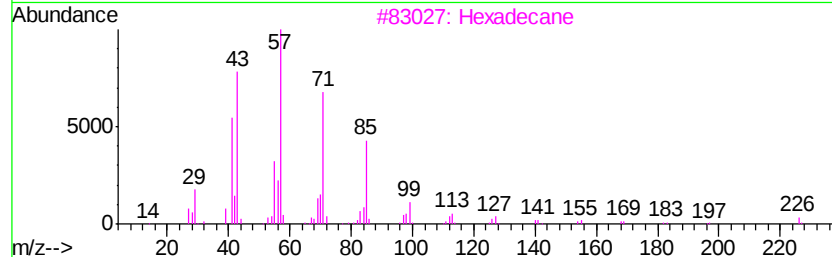
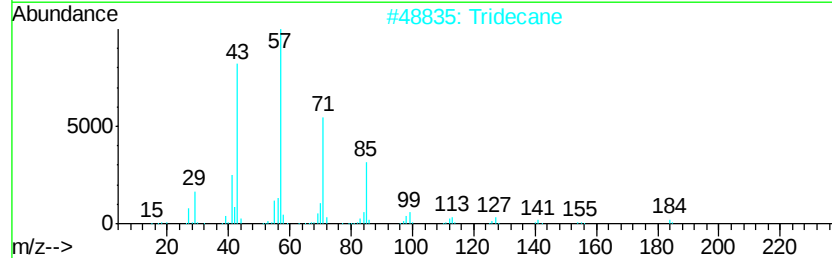
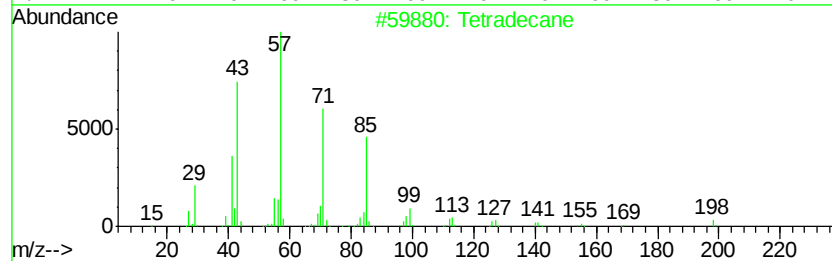
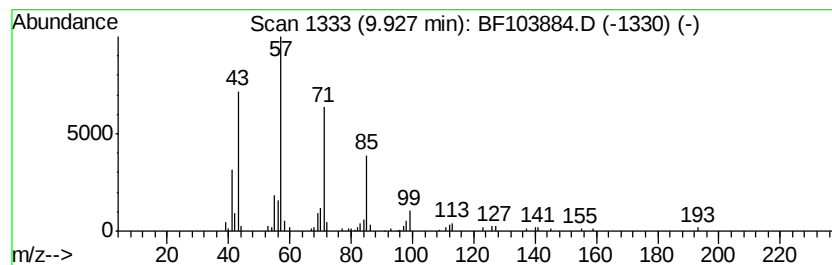
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Tetradecane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.93	2.18 ng	124889	Acenaphthene-d10	10.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetradecane	198	C14H30	000629-59-4	91
2		Tridecane	184	C13H28	000629-50-5	90
3		Hexadecane	226	C16H34	000544-76-3	87
4		1-Octadecanesulphonyl chloride	352	C18H37ClO2S	1000342-70-4	86
5		Decane, 2,3,5-trimethyl-	184	C13H28	062238-11-3	83



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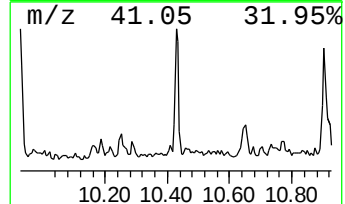
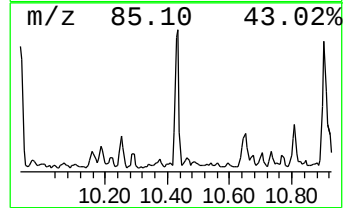
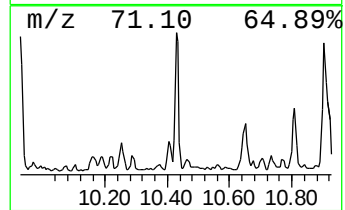
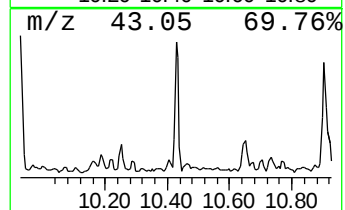
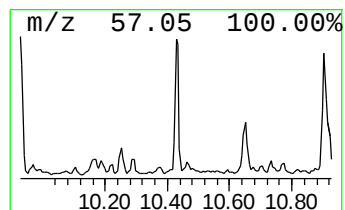
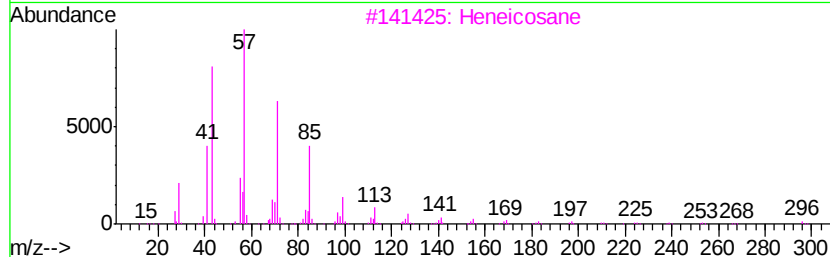
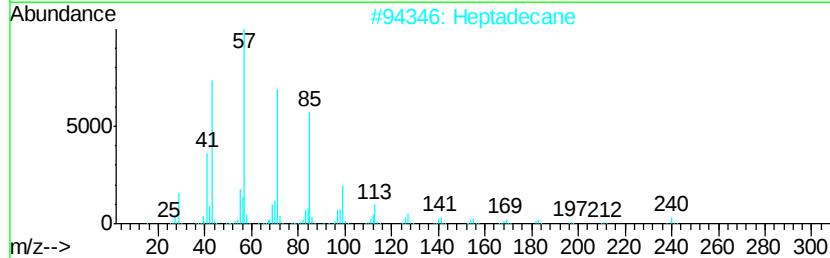
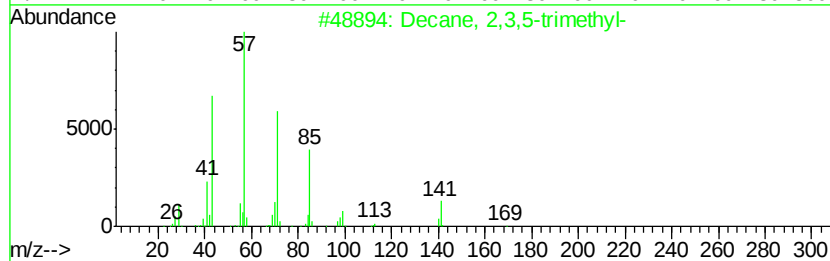
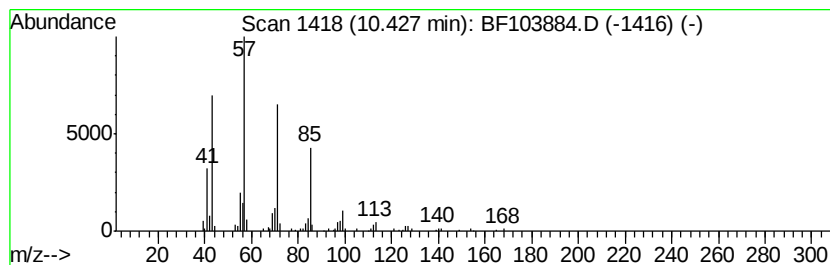
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 6 Decane, 2,3,5-trimethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.43	2.20 ng	126127	Acenaphthene-d10	10.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Decane, 2,3,5-trimethyl-	184	C13H28	062238-11-3	90
2		Heptadecane	240	C17H36	000629-78-7	90
3		Heneicosane	296	C21H44	000629-94-7	86
4		Tetradecane	198	C14H30	000629-59-4	86
5		Nonadecane	268	C19H40	000629-92-5	86



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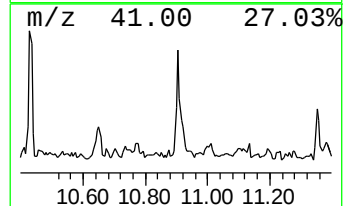
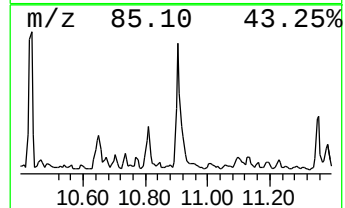
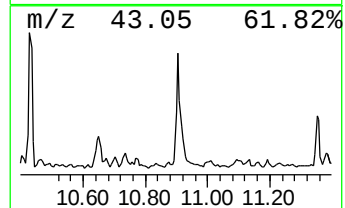
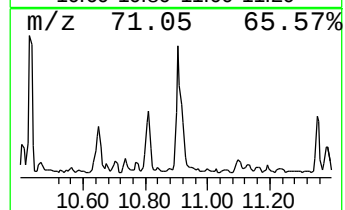
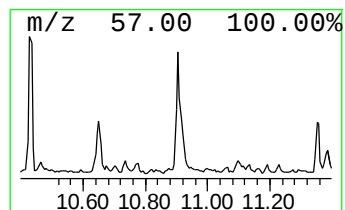
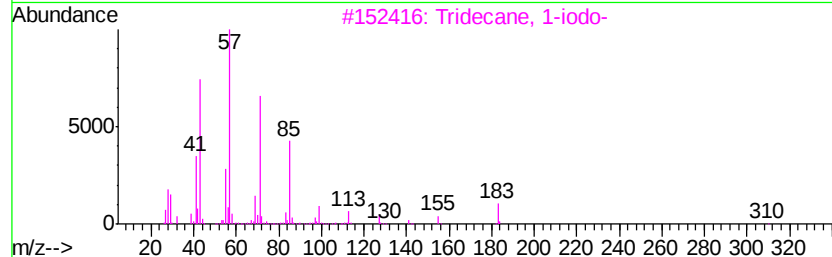
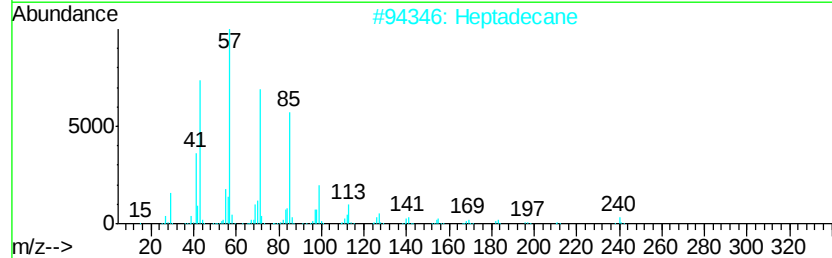
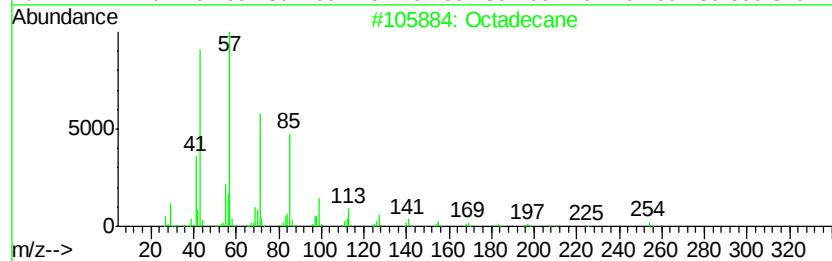
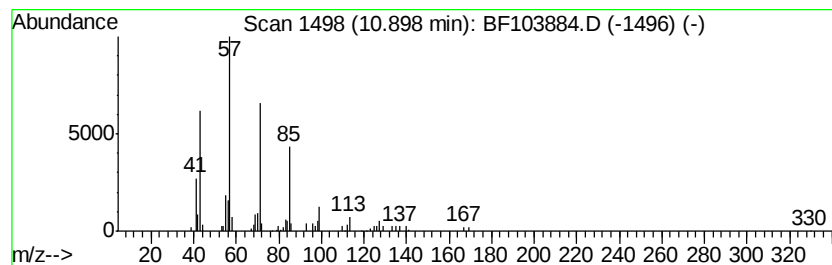
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Octadecane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.90	3.14 ng	159101	Phenanthrene-d10	11.51

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecane		254	C18H38	000593-45-3	91
2	Heptadecane		240	C17H36	000629-78-7	91
3	Tridecane, 1-iodo-		310	C13H27I	035599-77-0	90
4	Hexadecane		226	C16H34	000544-76-3	90
5	Pentadecane		212	C15H32	000629-62-9	87



Data Path : Z:\HPCHEM1\BNA F\DATA\BF032318\
Data File : BF103884.D
Acq On : 23 Mar 2018 17:00
Operator : JU/SJ
Sample : J1748-09
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
AU-05-032218-B

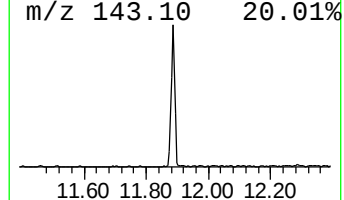
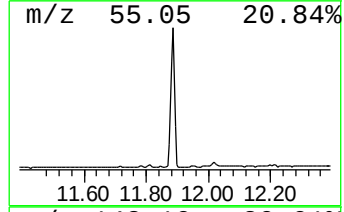
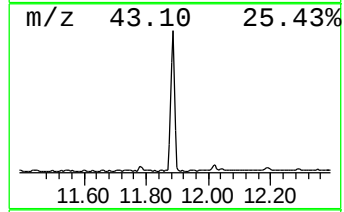
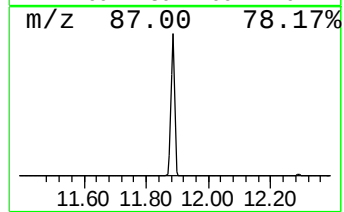
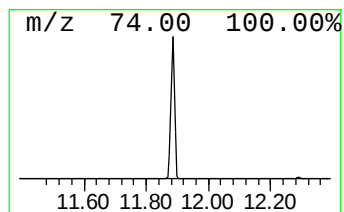
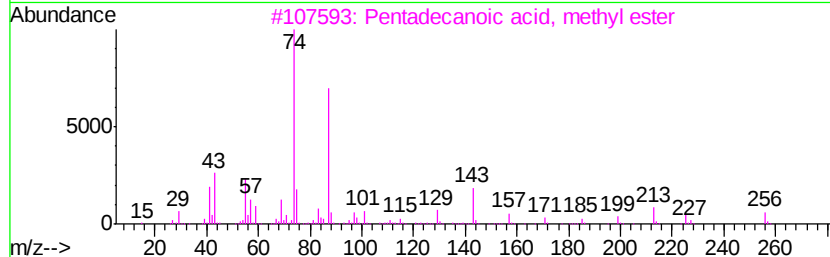
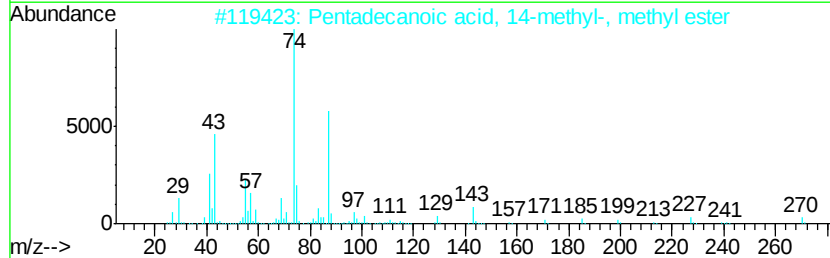
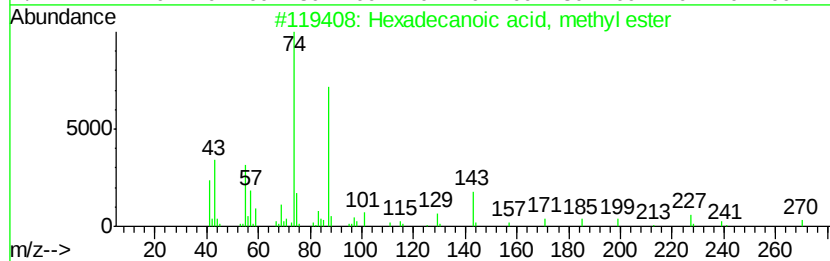
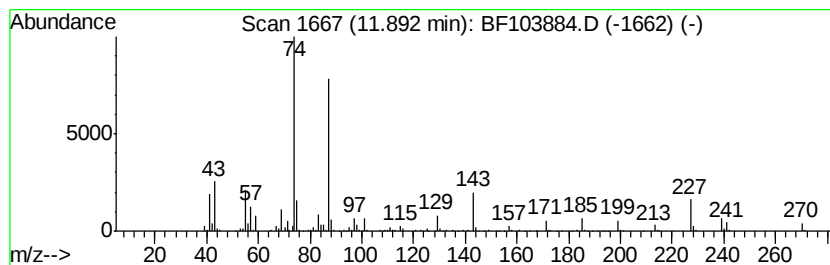
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Hexadecanoic acid, methyl e... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.89	29.46 ng	1490590	Phenanthrene-d10	11.51

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	93
3		Pentadecanoic acid, methyl ester	256	C16H32O2	007132-64-1	86
4		Hexadecanoic acid, 15-methyl-, m...	284	C18H36O2	006929-04-0	86
5		Methyl tetradecanoate	242	C15H30O2	000124-10-7	80



Data Path : Z:\HPCHEM1\BNA F\DATA\BF032318\
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Sample : J1748-09
Misc :
ALS Vial : 20 Sample Multiplier: 1

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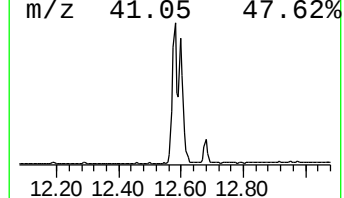
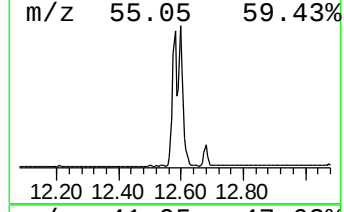
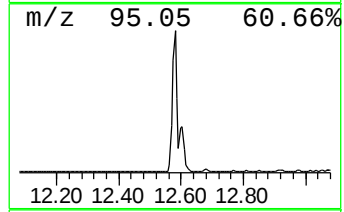
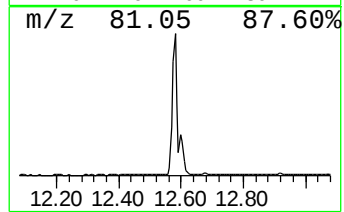
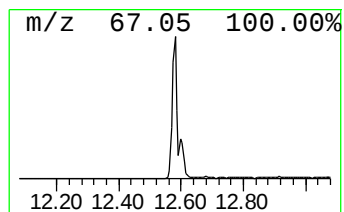
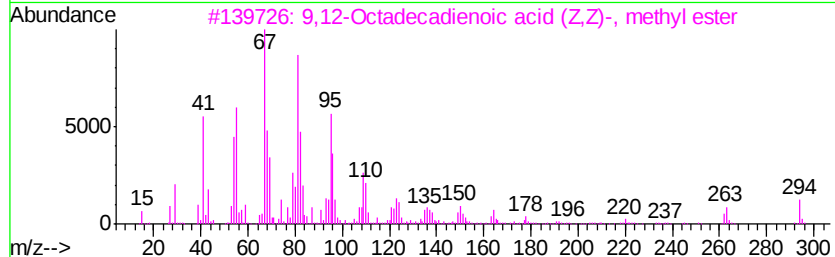
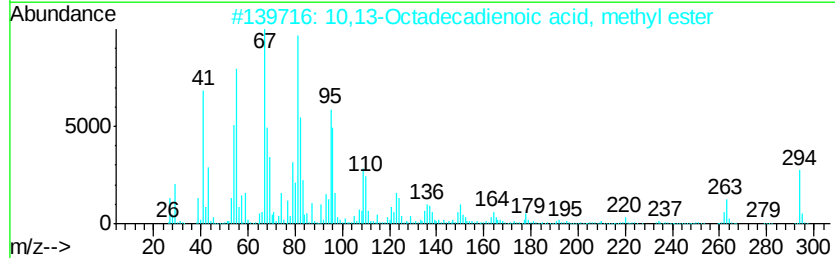
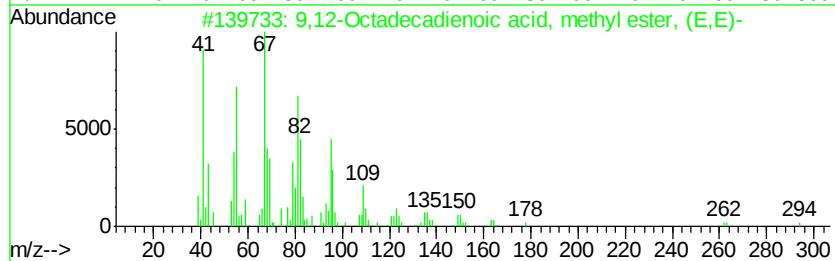
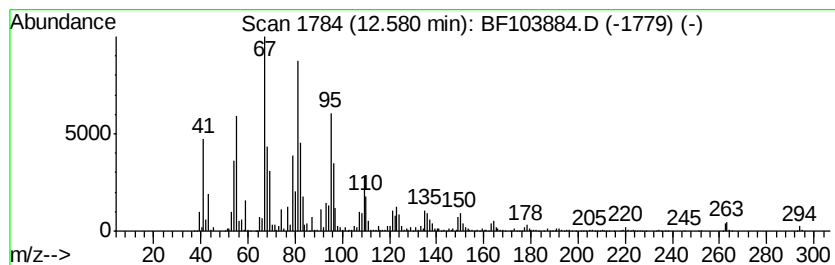
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.58	92.57 ng	4683620	Phenanthrene-d10	11.51

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9,12-Octadecadienoic acid, methy...	294	C19H34O2	002566-97-4	99
2		10,13-Octadecadienoic acid, meth...	294	C19H34O2	056554-62-2	99
3		9,12-Octadecadienoic acid (Z,Z)-...	294	C19H34O2	000112-63-0	99
4		Methyl 10-trans,12-cis-octadecad...	294	C19H34O2	1000336-44-2	98
5		11,14-Octadecadienoic acid, meth...	294	C19H34O2	056554-61-1	97



Data Path : Z:\HPCHEM1\BNA F\DATA\BF032318\
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Sample : J1748-09
Misc :
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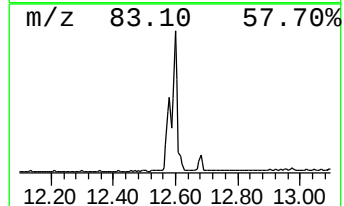
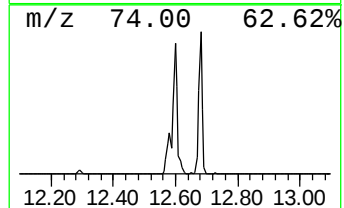
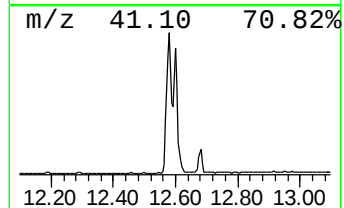
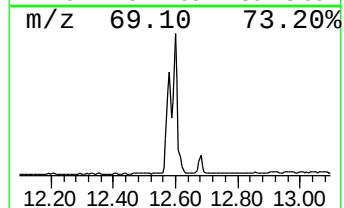
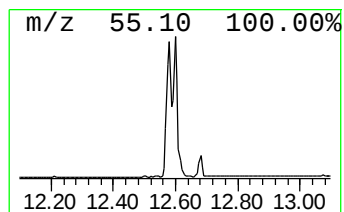
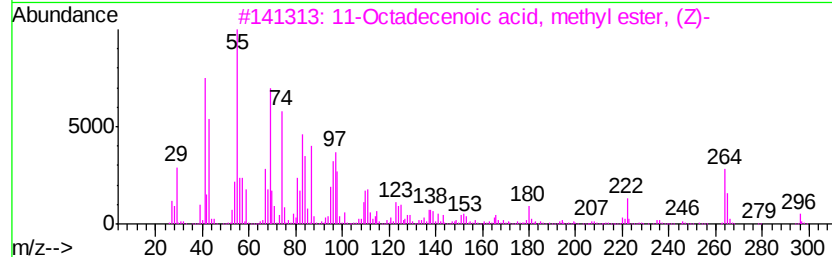
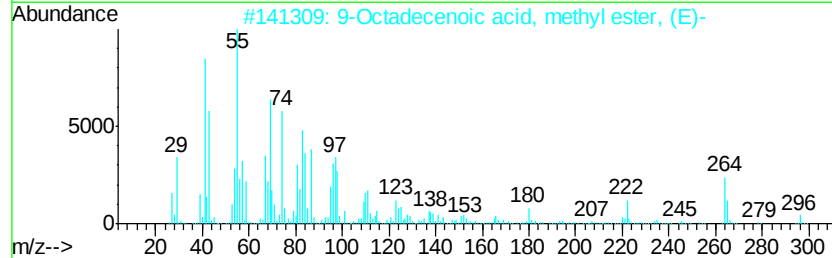
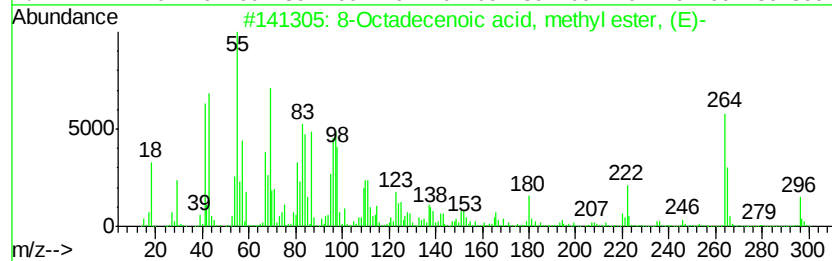
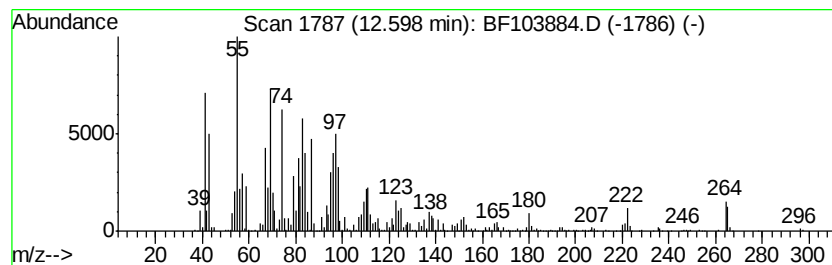
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 8-Octadecenoic acid, methyl... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.60	52.09 ng	2635650	Phenanthrene-d10	11.51

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		8-Octadecenoic acid, methyl este...	296	C19H36O2	026528-50-7	99
2		9-Octadecenoic acid, methyl este...	296	C19H36O2	001937-62-8	99
3		11-Octadecenoic acid, methyl est...	296	C19H36O2	001937-63-9	99
4		9-Octadecenoic acid, methyl ester	296	C19H36O2	002462-84-2	99
5		15-Octadecenoic acid, methyl ester	296	C19H36O2	004764-72-1	99



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Data File : BF103884.D
Acq On : 23 Mar 2018 17:00
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Sample : J1748-09
Misc :
ALS Vial : 20 Sample Multiplier: 1

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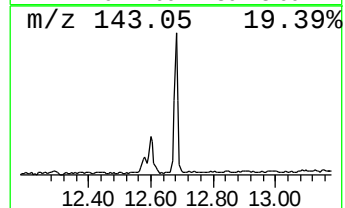
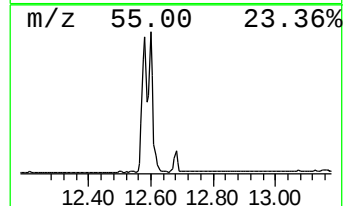
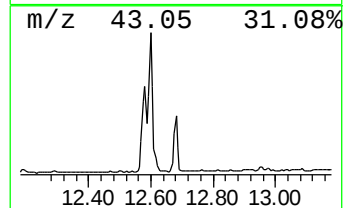
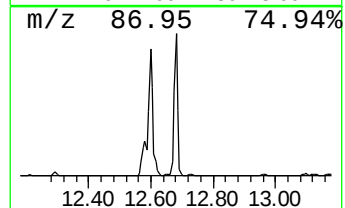
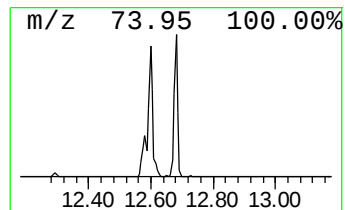
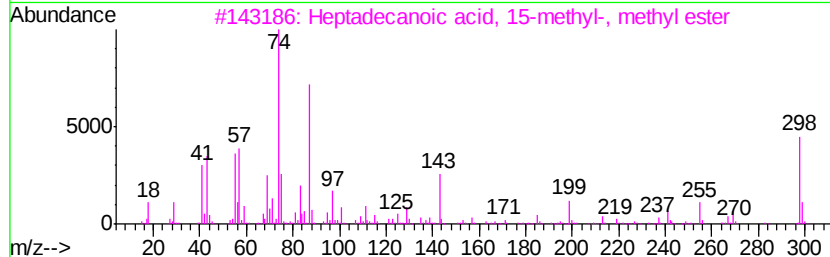
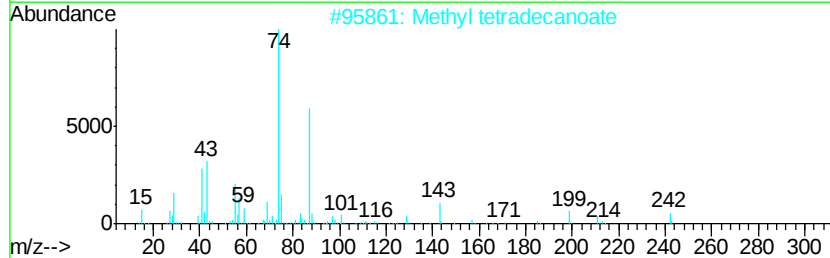
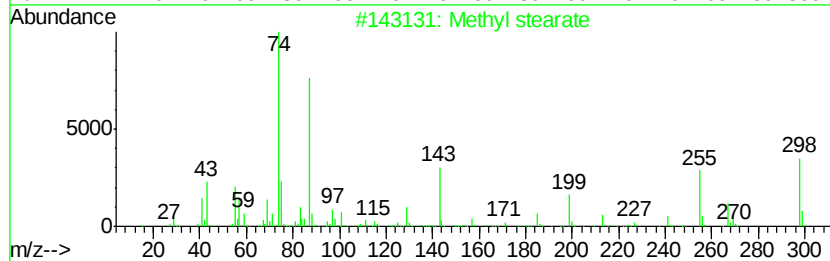
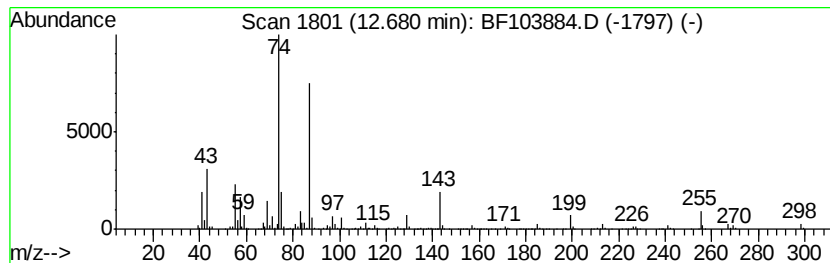
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Methyl stearate Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.68	12.11 ng	612580	Phenanthrene-d10	11.51

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Methyl stearate	298	C19H38O2	000112-61-8	98
2		Methyl tetradecanoate	242	C15H30O2	000124-10-7	95
3		Heptadecanoic acid, 15-methyl-, ...	298	C19H38O2	054833-55-5	91
4		Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	90
5		Heptadecanoic acid, 16-methyl-, ...	298	C19H38O2	005129-61-3	90



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ALS Vial : 20 Sample Multiplier: 1

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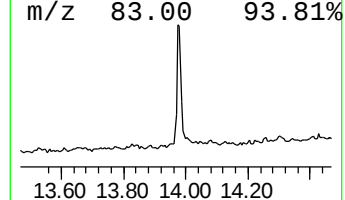
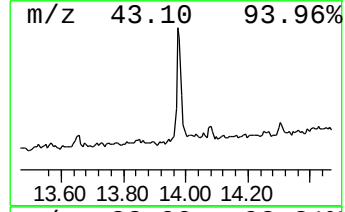
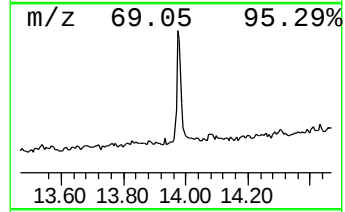
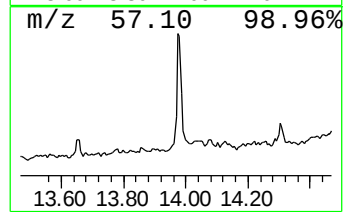
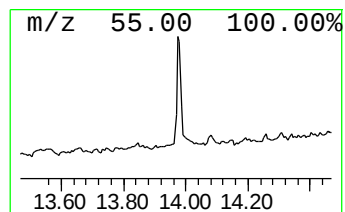
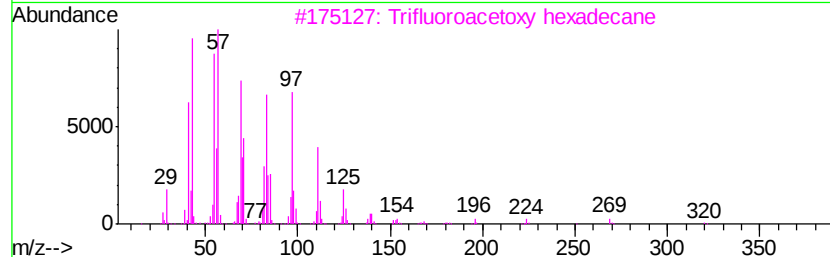
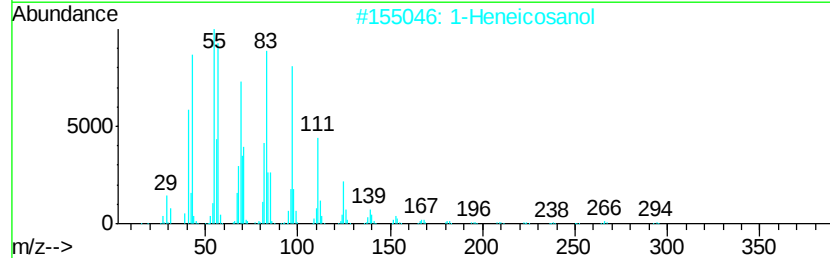
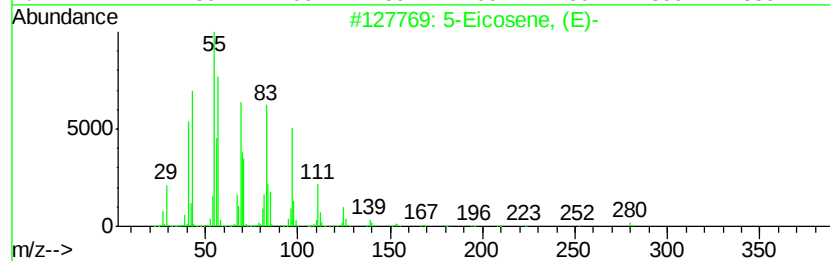
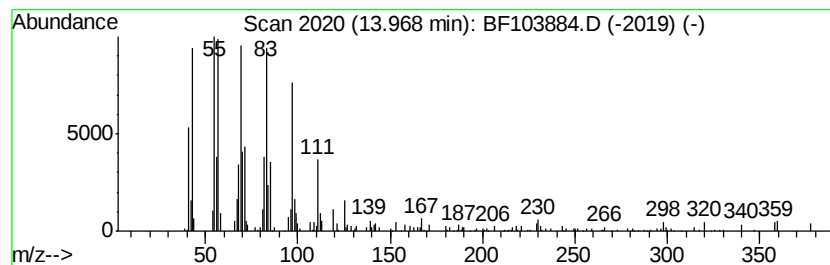
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 5-Eicosene, (E)- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.97	8.77 ng	354693	Chrysene-d12	14.17

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			5-Eicosene, (E)-	280	C20H40	074685-30-6	97
2			1-Heneicosanol	312	C21H44O	015594-90-8	94
3			Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	94
4			Behenic alcohol	326	C22H46O	000661-19-8	94
5			Trifluoroacetic acid, pentadecyl...	324	C17H31F3O2	959010-23-2	91



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032318\
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 Sample : J1748-09
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031518.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...	2.30	3.4	ng	146558	1	6.97	873653	20.0
2-Pentanone, 4-hy...	5.22	3.4	ng	147216	1	6.97	873653	20.0
unknown6.75	6.75	69.1	ng	3019710	1	6.97	873653	20.0
Tetradecane	9.93	2.2	ng	124889	3	10.02	1145260	20.0
Decane, 2,3,5-tri...	10.43	2.2	ng	126127	3	10.02	1145260	20.0
Octadecane	10.90	3.1	ng	159101	4	11.51	1011880	20.0
Hexadecanoic acid...	11.89	29.5	ng	1490590	4	11.51	1011880	20.0
9,12-Octadecadien...	12.58	92.6	ng	4683620	4	11.51	1011880	20.0
8-Octadecenoic ac...	12.60	52.1	ng	2635650	4	11.51	1011880	20.0
Methyl stearate	12.68	12.1	ng	612580	4	11.51	1011880	20.0
5-Eicosene, (E)-	13.97	8.8	ng	354693	5	14.17	808773	20.0