

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF112215\
 Data File : BF083181.D
 Acq On : 23 Nov 2015 2:04
 Operator : UM/IZ
 Sample : G4512-02
 Misc :
 ALS Vial : 67 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FILL-DIRT-2

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-BF112115.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	3.427	119	126	135	rBV	5673455	14016476	100.00%	18.295%
2	4.799	243	246	262	rVB	1900976	3550894	25.33%	4.635%
3	5.256	284	286	289	rBV	4957484	5738648	40.94%	7.490%
4	6.319	376	379	381	rBV	5600657	6270130	44.73%	8.184%
5	6.422	386	388	391	rBV	5099458	5668182	40.44%	7.398%
6	6.650	405	408	410	rBV	1468005	1395432	9.96%	1.821%
7	6.810	419	422	424	rBV	4131252	4639347	33.10%	6.056%
8	7.222	455	458	460	rBV	3930606	3961695	28.26%	5.171%
9	7.942	518	521	525	rBV2	1375864	2005193	14.31%	2.617%
10	9.016	612	615	618	rBV	5557863	5671185	40.46%	7.402%
11	9.416	648	650	653	rBV	1162360	1017143	7.26%	1.328%
12	9.633	667	669	671	rBV	161370	144507	1.03%	0.189%
13	9.679	671	673	675	rVV	1280759	1808510	12.90%	2.361%
14	10.136	705	713	714	rBV	308869	333234	2.38%	0.435%
15	10.354	729	732	734	rBV	127422	203454	1.45%	0.266%
16	10.468	740	742	745	rBV2	2697905	3188779	22.75%	4.162%
17	10.605	752	754	758	rBV	342968	399410	2.85%	0.521%
18	11.051	791	793	795	rBV3	180070	162672	1.16%	0.212%
19	11.165	800	803	806	rBV	1250364	1957738	13.97%	2.555%
20	11.234	806	809	812	rVB2	141481	177794	1.27%	0.232%
21	11.714	849	851	854	rVV2	974875	1002423	7.15%	1.308%
22	11.771	854	856	861	rVB2	104474	189221	1.35%	0.247%
23	12.365	906	908	910	rBV	687218	668106	4.77%	0.872%
24	12.491	917	919	922	rVV2	352200	431639	3.08%	0.563%
25	12.582	925	927	930	rVV	476305	713847	5.09%	0.932%
26	12.742	938	941	944	rVV	3745010	4191499	29.90%	5.471%
27	12.925	953	957	959	rBV2	75567	177252	1.26%	0.231%
28	13.222	981	983	987	rVB2	131380	154597	1.10%	0.202%
29	13.657	1018	1021	1024	rBV2	532511	715810	5.11%	0.934%
30	13.782	1029	1032	1038	rVV2	1558989	2084456	14.87%	2.721%
31	13.885	1038	1041	1043	rBV2	75911	157933	1.13%	0.206%
32	14.285	1073	1076	1082	rVV4	176182	423047	3.02%	0.552%
33	14.777	1117	1119	1123	rBV	305069	554405	3.96%	0.724%
34	14.868	1125	1127	1128	rBV	151008	158623	1.13%	0.207%

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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-BF112115.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	15.040	1140	1142	1145	rVB	192522	237041	1.69%	0.309%
36	15.097	1145	1147	1149	rBV	233457	302877	2.16%	0.395%
37	15.154	1149	1152	1154	rBV	818023	1061588	7.57%	1.386%
38	15.577	1186	1189	1191	rBV	163021	279952	2.00%	0.365%
39	15.623	1191	1193	1196	rVB3	95867	187174	1.34%	0.244%
40	16.411	1259	1262	1267	rVB2	103746	213892	1.53%	0.279%
41	16.548	1272	1274	1278	rBV3	119631	240623	1.72%	0.314%
42	16.788	1292	1295	1301	rBV	77365	156361	1.12%	0.204%

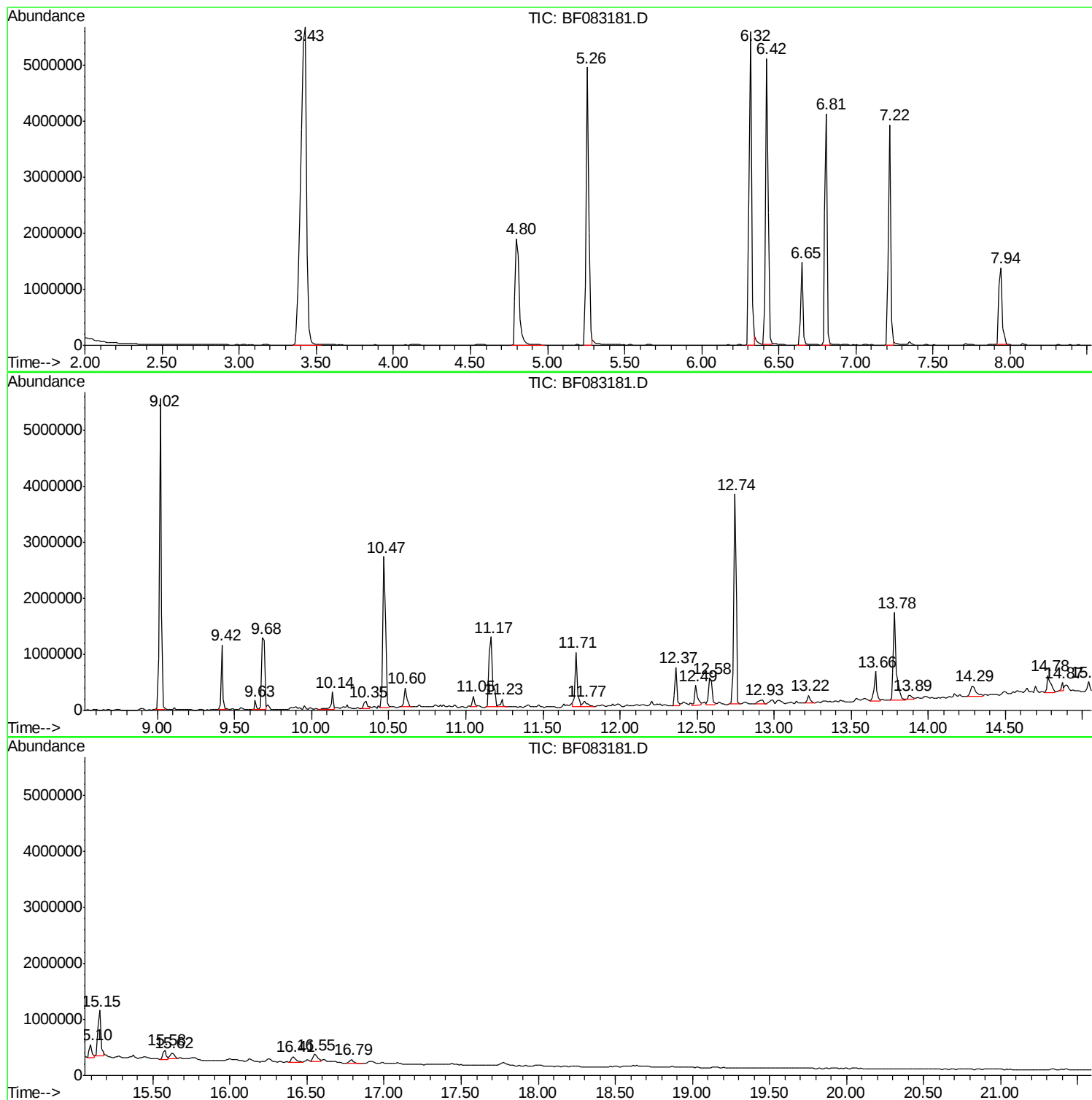
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF112115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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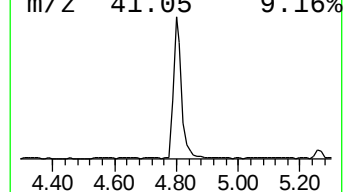
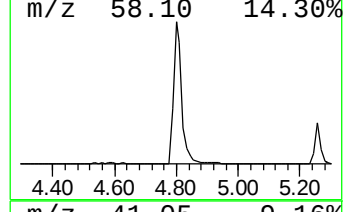
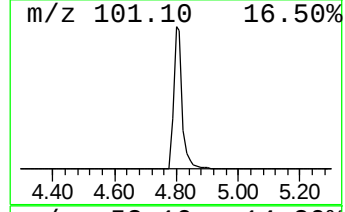
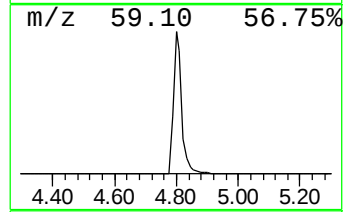
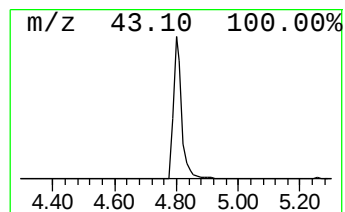
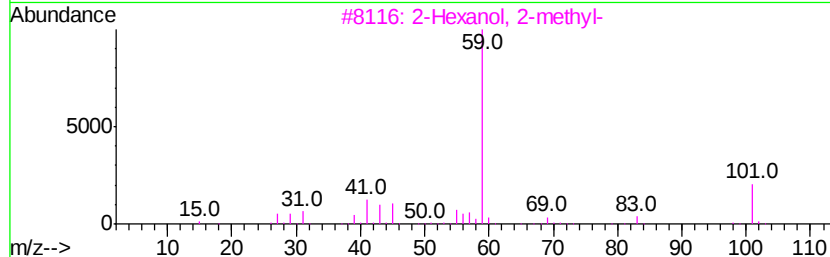
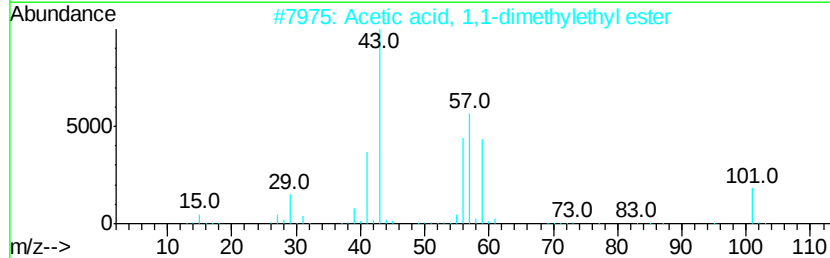
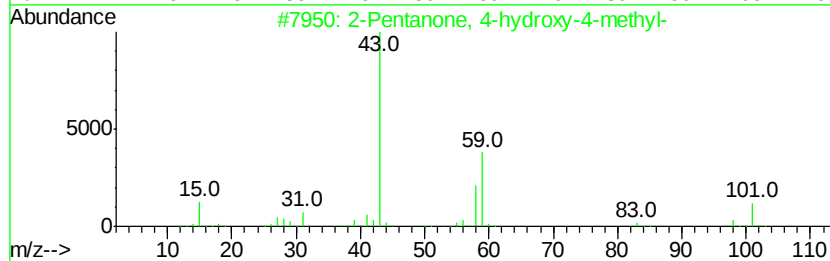
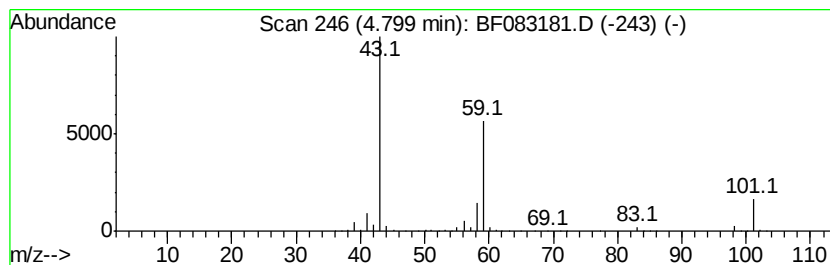
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF112115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown4.80 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.80	50.89 ng	3550890	1,4-Dichlorobenzene-d4	6.65

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	40
2		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	33
4		3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
5		Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	32



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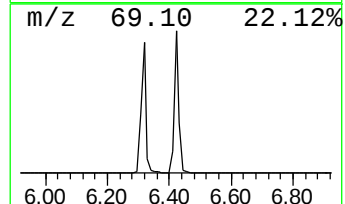
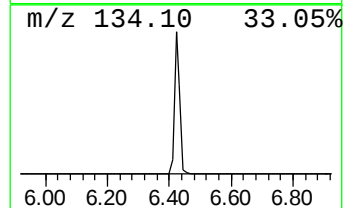
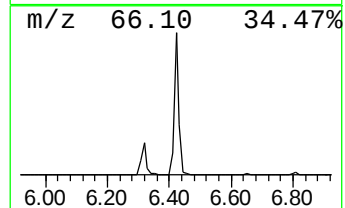
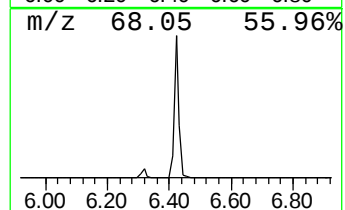
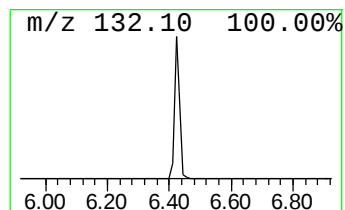
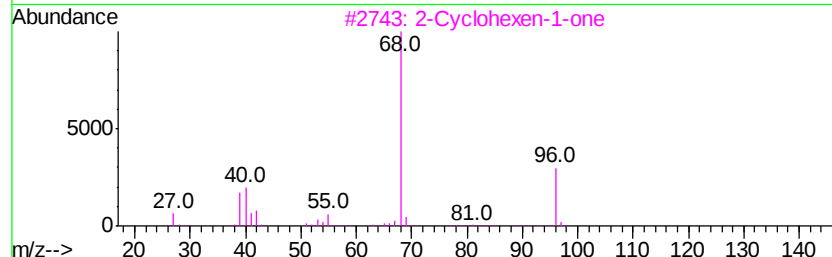
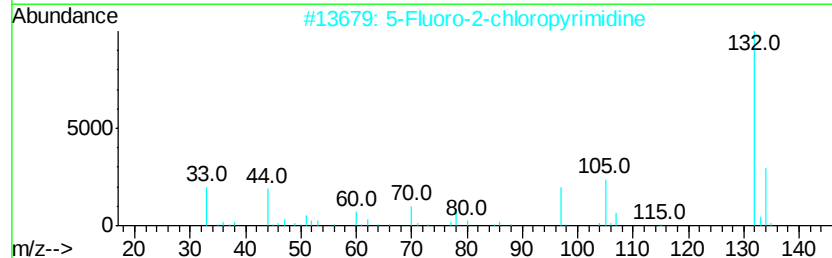
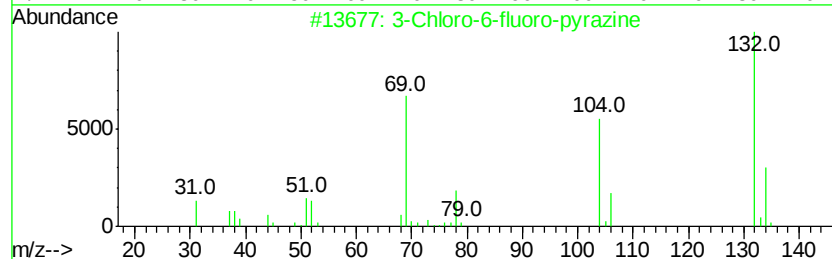
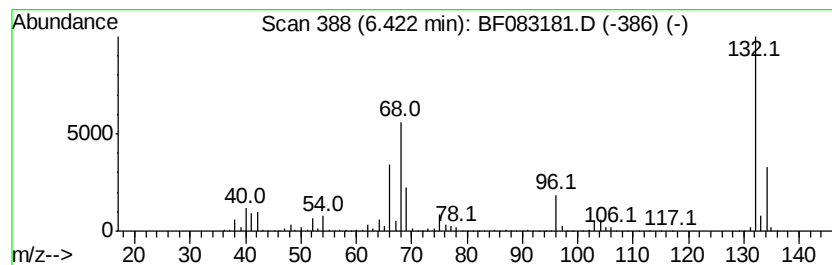
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 3 unknown6.42 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.42	81.24 ng	5668180	1,4-Dichlorobenzene-d4	6.65

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	12
3		2-Cyclohexen-1-one	96	C6H8O	000930-68-7	10
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		(5-METHYL-2-PYRIDYL)ACETONITRILE	132	C8H8N2	1000241-93-9	10



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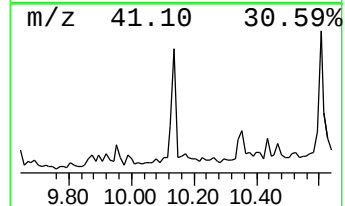
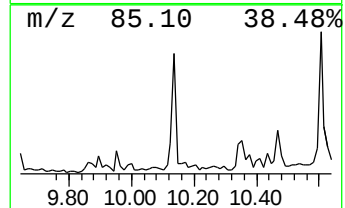
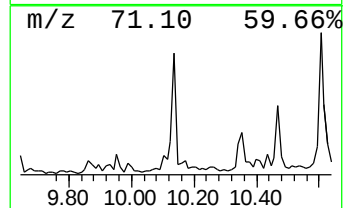
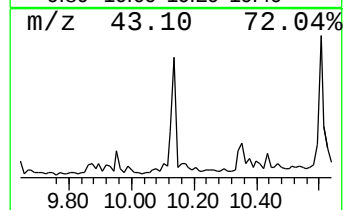
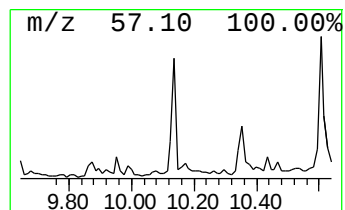
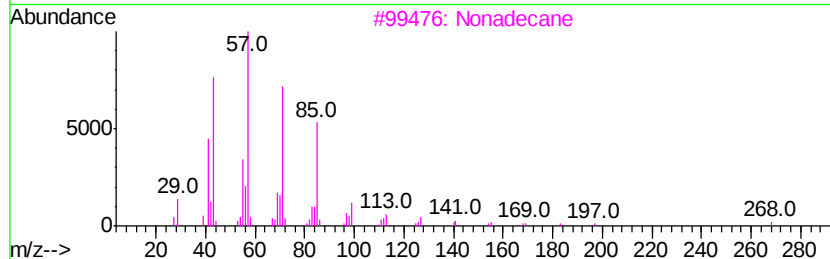
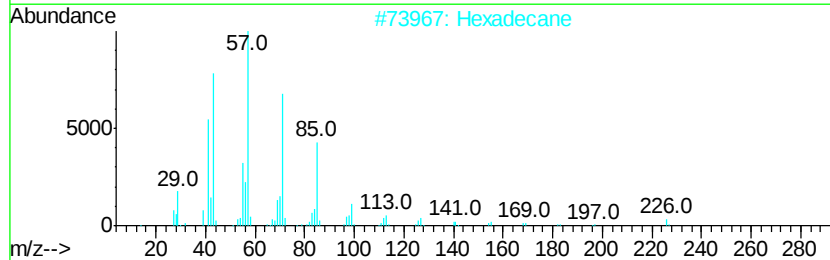
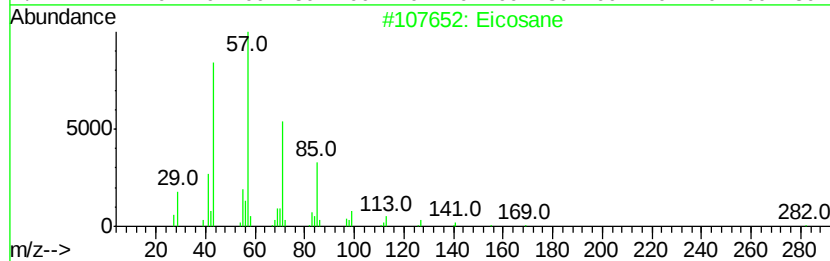
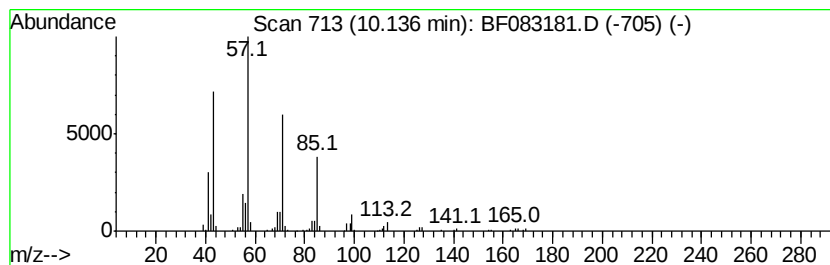
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Eicosane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.14	3.69 ng	333234	Acenaphthene-d10	9.68

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane		282	C20H42	000112-95-8	91
2	Hexadecane		226	C16H34	000544-76-3	87
3	Nonadecane		268	C19H40	000629-92-5	86
4	Tetracosane		338	C24H50	000646-31-1	83
5	Eicosane, 10-methyl-		296	C21H44	054833-23-7	83



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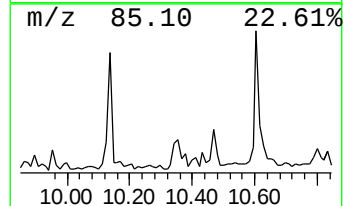
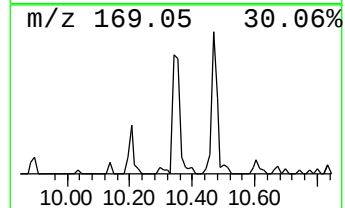
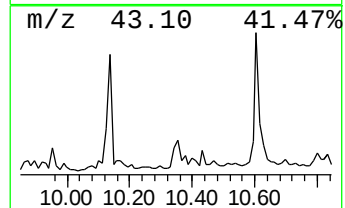
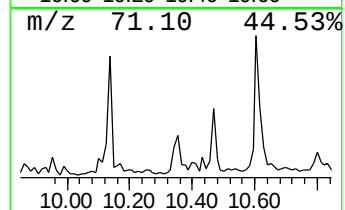
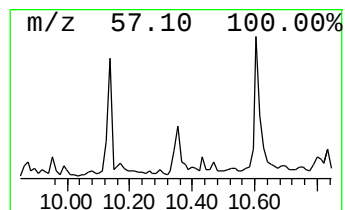
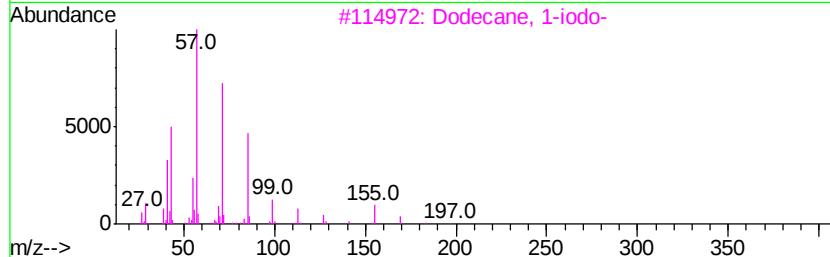
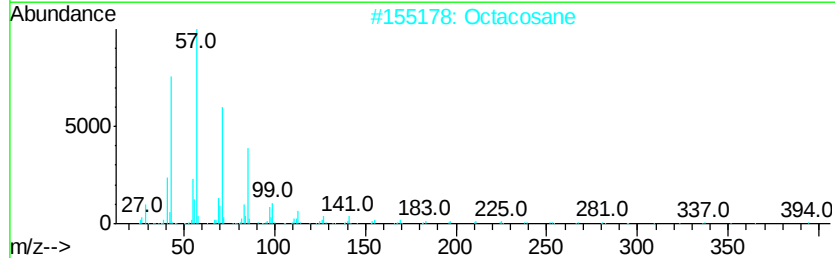
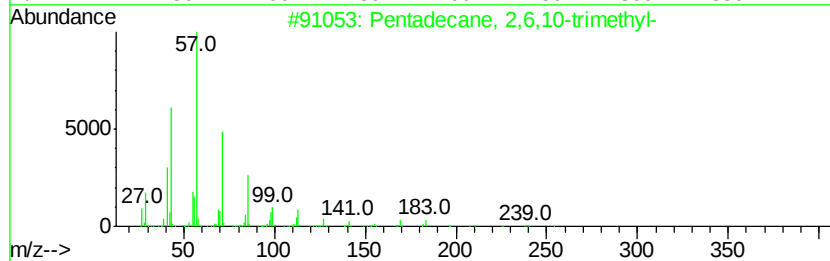
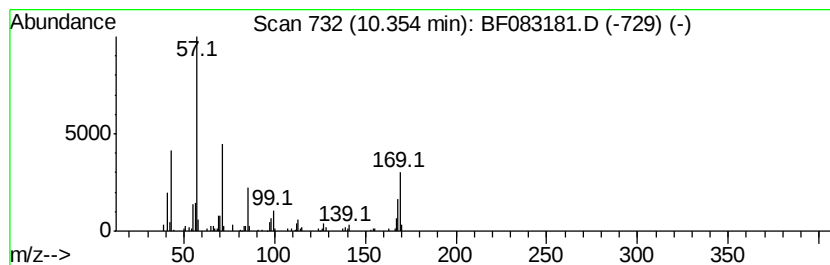
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Pentadecane, 2,6,10-trimethyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.35	2.25 ng	203454	Acenaphthene-d10	9.68

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentadecane, 2,6,10-trimethyl-	254	C18H38	003892-00-0	53
2		Octacosane	394	C28H58	000630-02-4	52
3		Dodecane, 1-iodo-	296	C12H25I	004292-19-7	50
4		Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	49
5		Tetracosane	338	C24H50	000646-31-1	49



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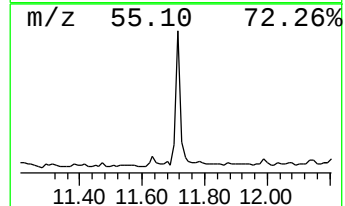
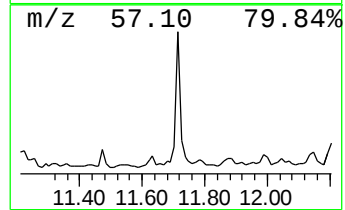
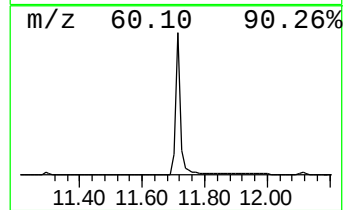
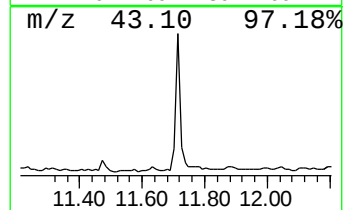
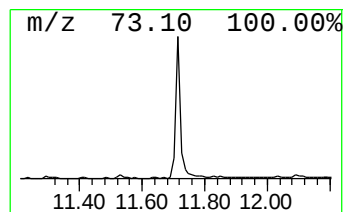
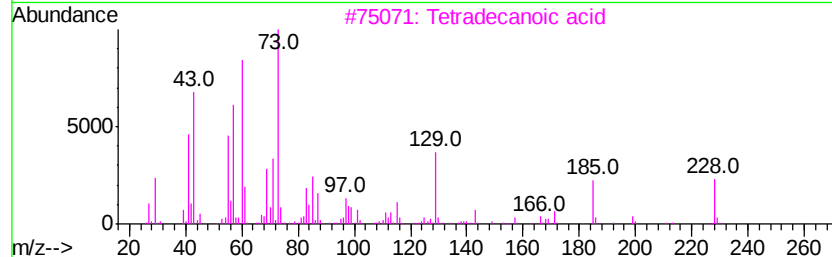
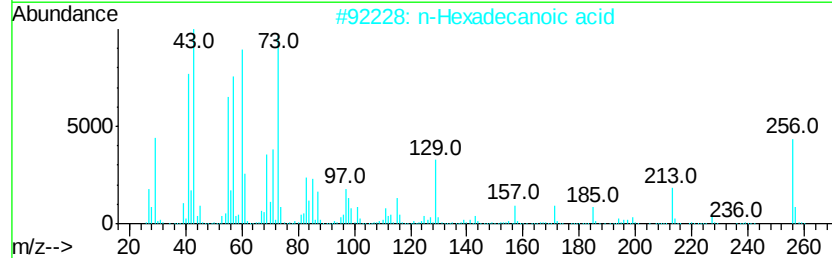
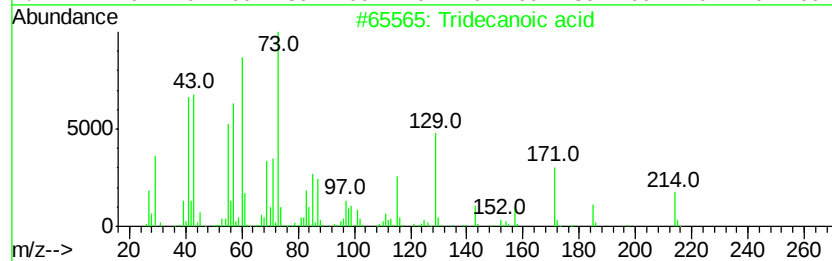
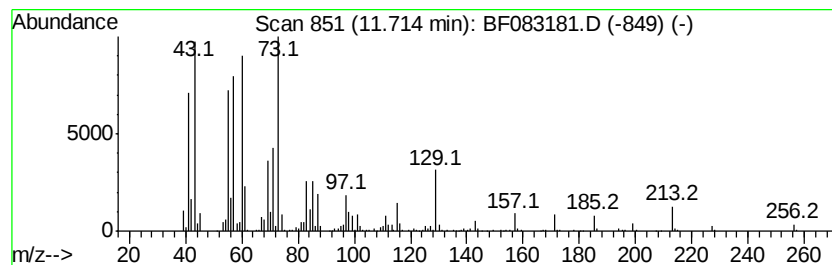
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Peak Number 8 Tridecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.71	10.24 ng	1002420	Phenanthrene-d10	11.16

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tridecanoic acid	214	C13H26O2	000638-53-9	95
2		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	93
3		Tetradecanoic acid	228	C14H28O2	000544-63-8	91
4		Pentadecanoic acid	242	C15H30O2	001002-84-2	87
5		Undecanoic acid	186	C11H22O2	000112-37-8	86



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF112215\
Data File : BF083181.D
Acq On : 23 Nov 2015 2:04
Operator : UM/IZ
Sample : G4512-02
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
FILL-DIRT-2

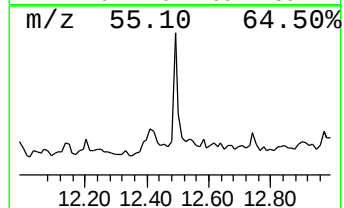
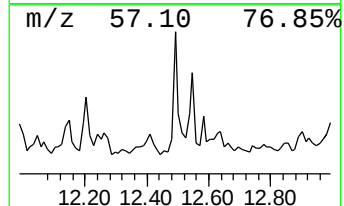
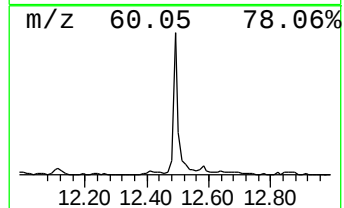
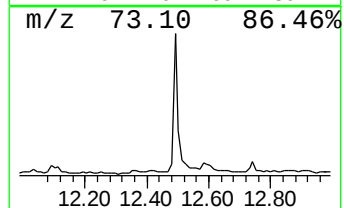
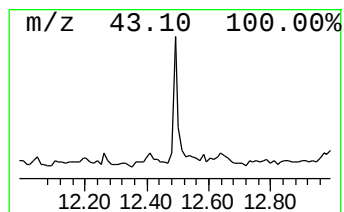
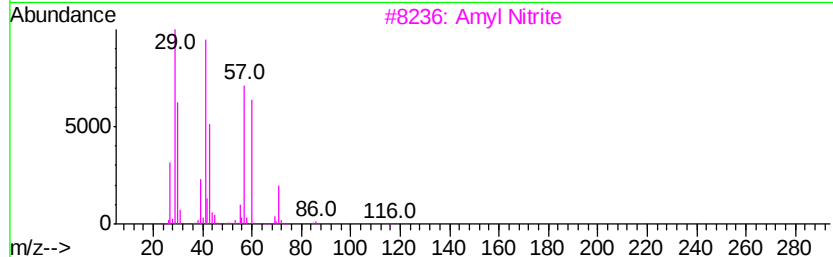
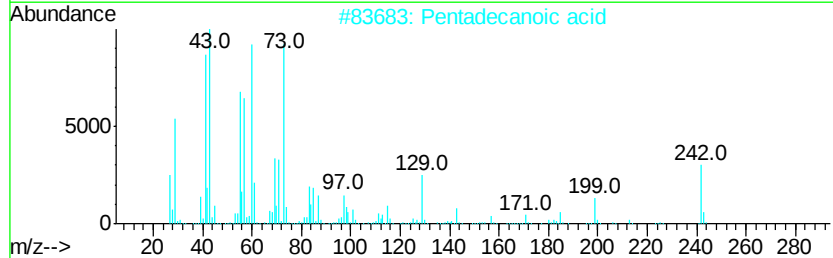
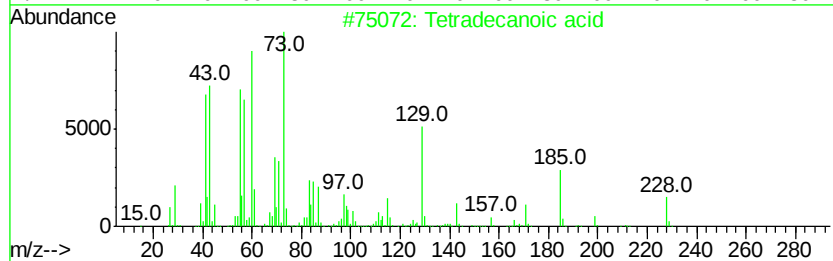
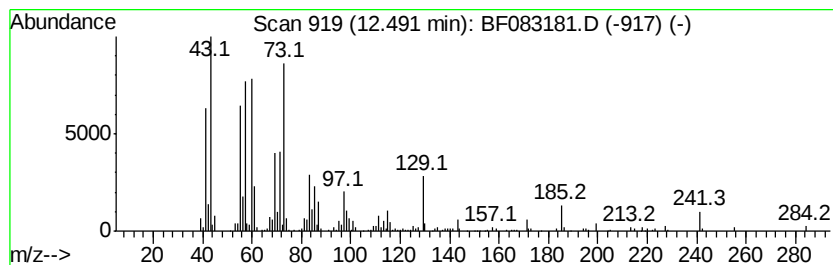
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Tetradecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.49	4.14 ng	431639	Chrysene-d12	13.78

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tetradecanoic acid	228	C14H28O2	000544-63-8	93
2		Pentadecanoic acid	242	C15H30O2	001002-84-2	58
3		Amvl Nitrite	117	C5H11NO2	000110-46-3	27
4		Undecane	156	C11H24	001120-21-4	11
5		Tetradecane, 1-iodo-	324	C14H29I	019218-94-1	11



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF112215\
Data File : BF083181.D
Acq On : 23 Nov 2015 2:04
Operator : UM/IZ
Sample : G4512-02
Misc :
ALS Vial : 67 Sample Multiplier: 1

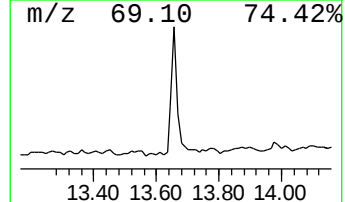
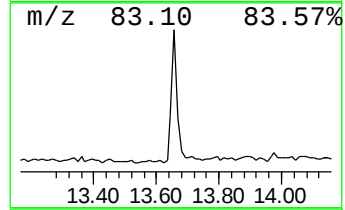
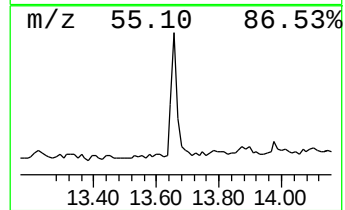
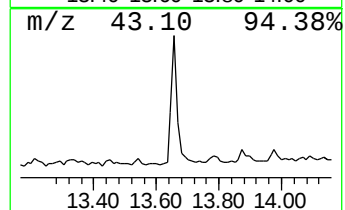
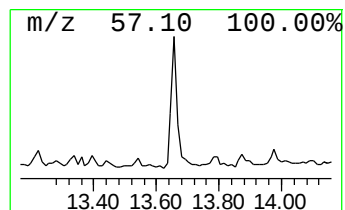
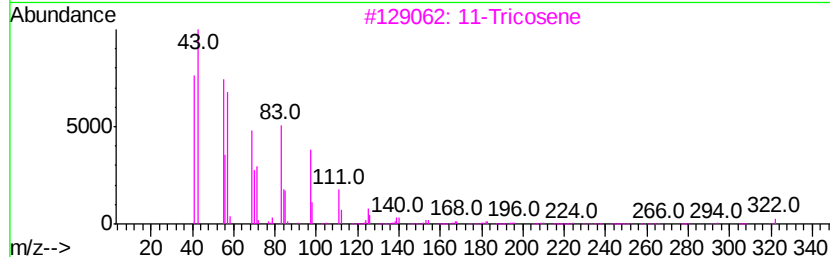
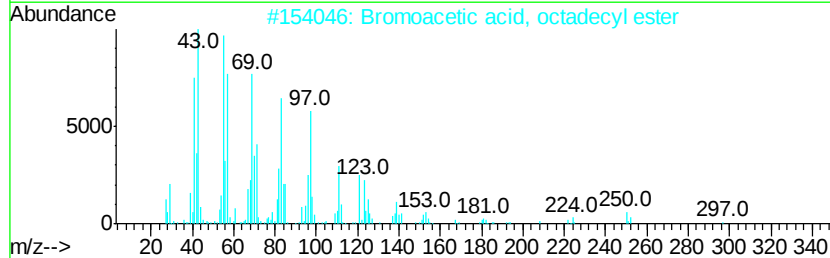
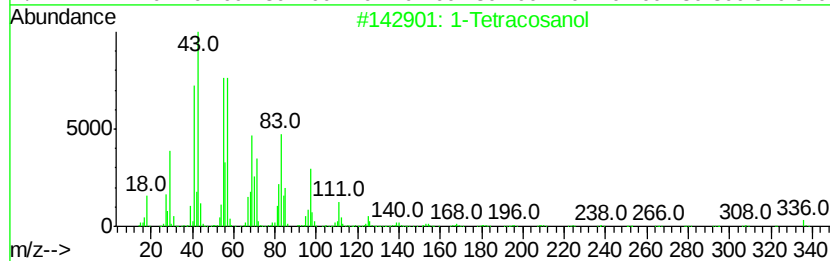
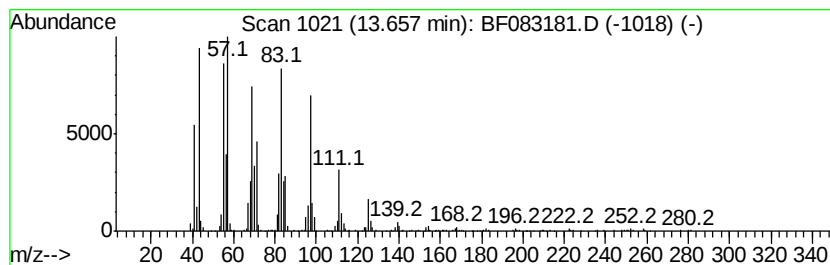
Instrument :
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ClientSampleId :
FILL-DIRT-2

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF112115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 1-Tetracosanol Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.66	6.87 ng	715810	Chrysene-d12	13.78	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Tetracosanol	354	C24H50O	000506-51-4	91
2	Bromoacetic acid, octadecyl ester	390	C20H39BrO2	018992-03-5	91
3	11-Tricosene	322	C23H46	052078-56-5	91
4	1-Octadecene	252	C18H36	000112-88-9	87
5	Bromoacetic acid, pentadecyl ester	348	C17H33BrO2	131143-01-6	87



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF112215\
Data File : BF083181.D
Acq On : 23 Nov 2015 2:04
Operator : UM/IZ
Sample : G4512-02
Misc :
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Instrument :
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ClientSampleId :
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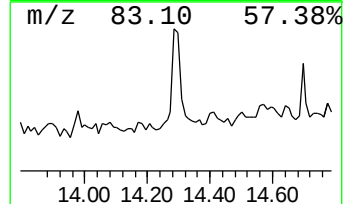
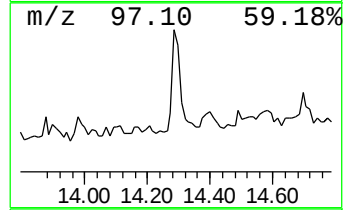
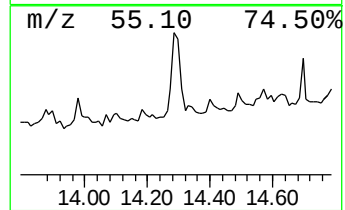
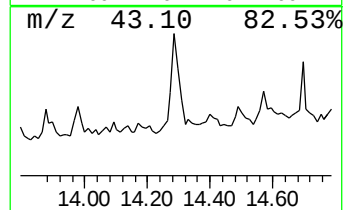
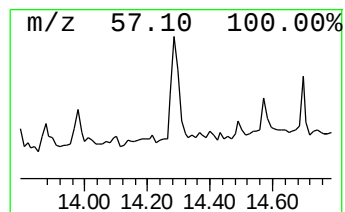
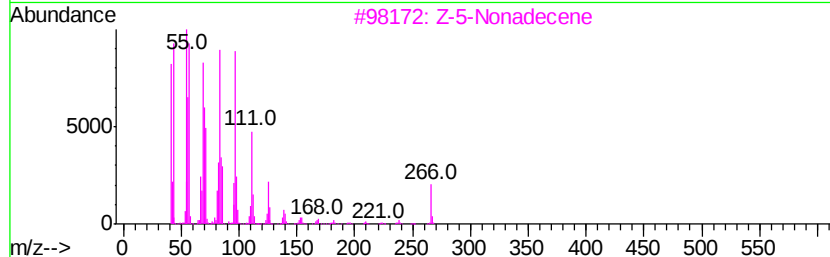
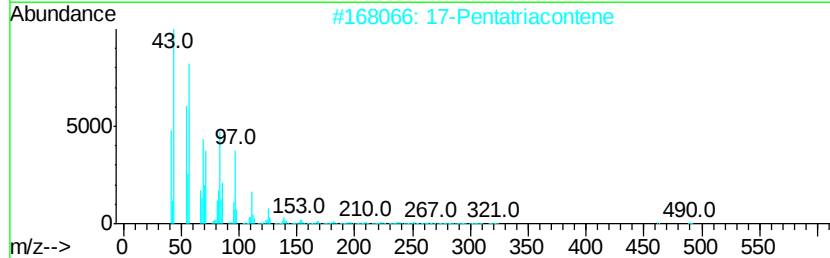
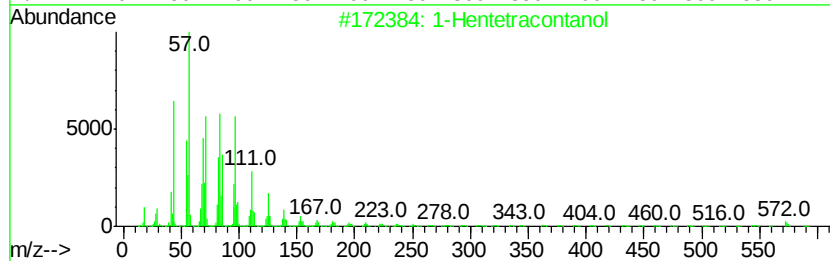
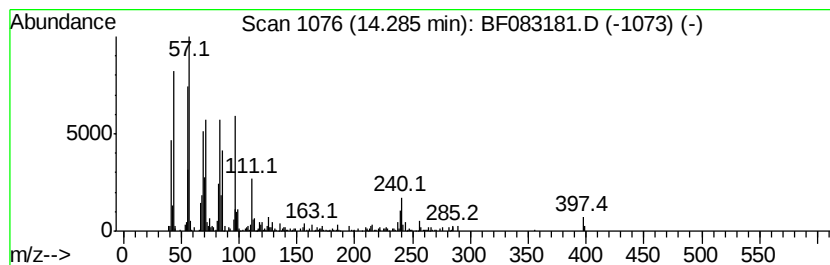
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 1-Hentetracontanol Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.29	4.06 ng	423047	Chrysene-d12	13.78

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Hentetracontanol	593	C41H84O	040710-42-7	74
2			17-Pentatriacontene	491	C35H70	006971-40-0	68
3			Z-5-Nonadecene	266	C19H38	1000131-11-8	53
4			1-Heptadecanol	256	C17H36O	001454-85-9	45
5			1-Hexacosanol	382	C26H54O	000506-52-5	45



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF112215\
Data File : BF083181.D
Acq On : 23 Nov 2015 2:04
Operator : UM/IZ
Sample : G4512-02
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument :
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ClientSampleId :
FILL-DIRT-2

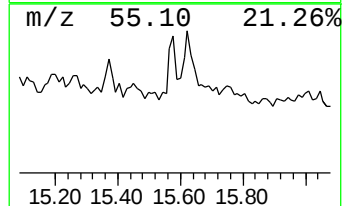
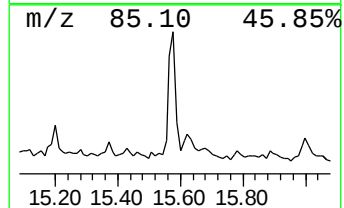
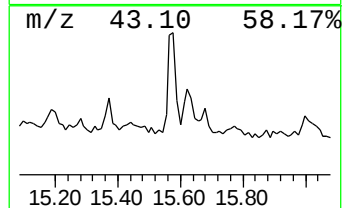
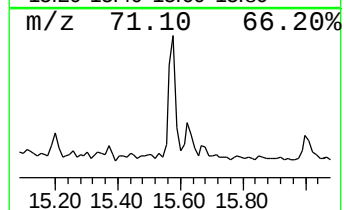
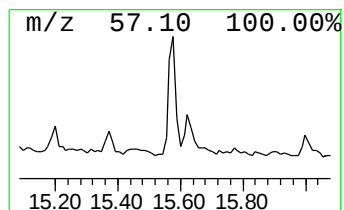
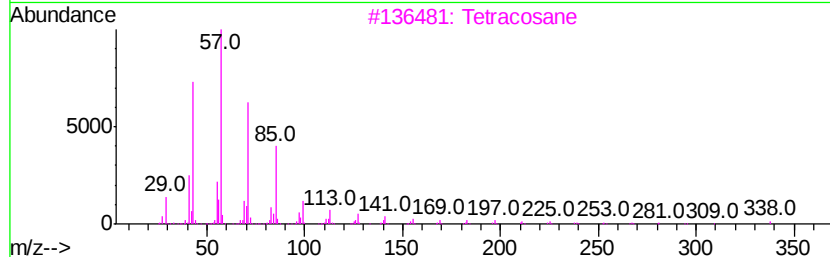
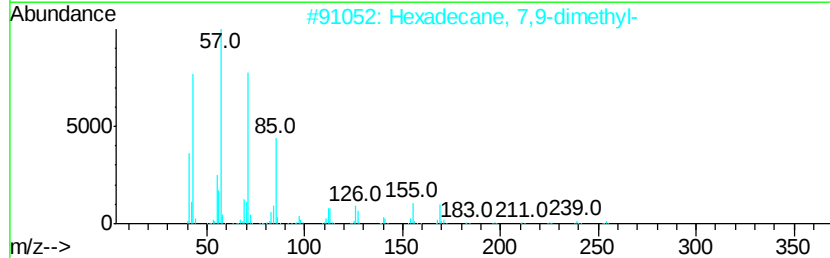
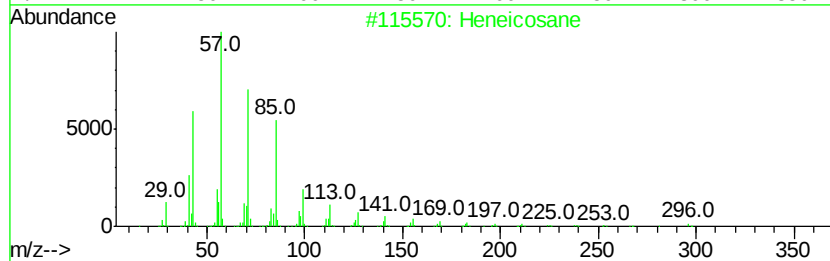
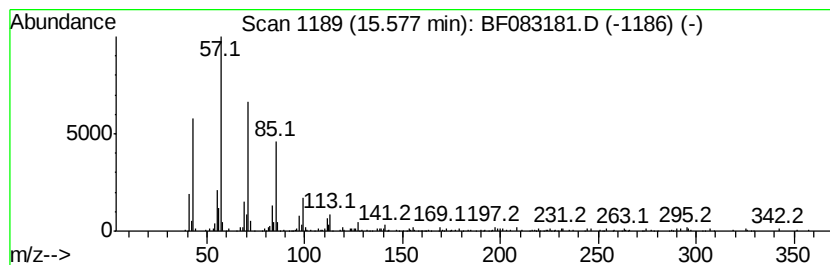
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 Heneicosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.58	5.27 ng	279952	Perylene-d12	15.15

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane		296	C21H44	000629-94-7	90
2	Hexadecane, 7,9-dimethyl-		254	C18H38	021164-95-4	90
3	Tetracosane		338	C24H50	000646-31-1	83
4	Dodecane, 2-methyl-6-propyl-		226	C16H34	055045-08-4	83
5	Tetradecane, 4-ethyl-		226	C16H34	055045-14-2	80



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF112215\
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Sample : G4512-02
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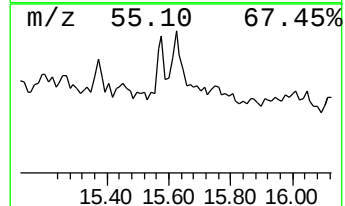
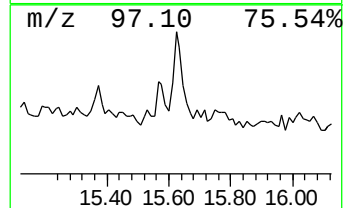
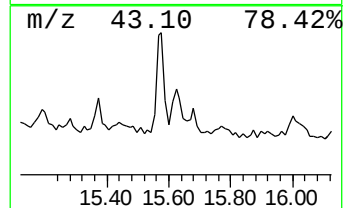
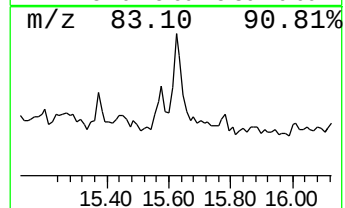
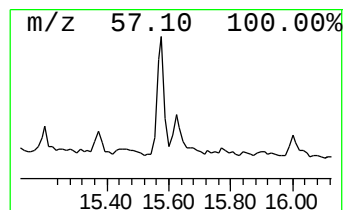
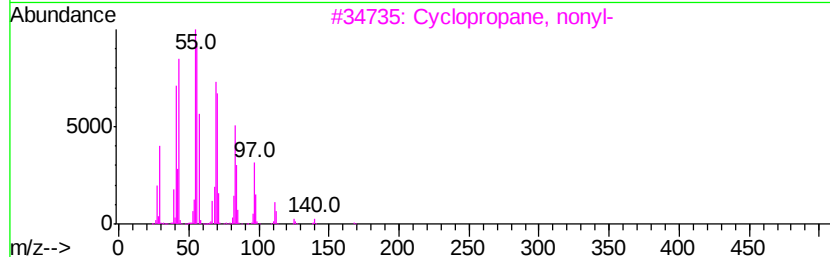
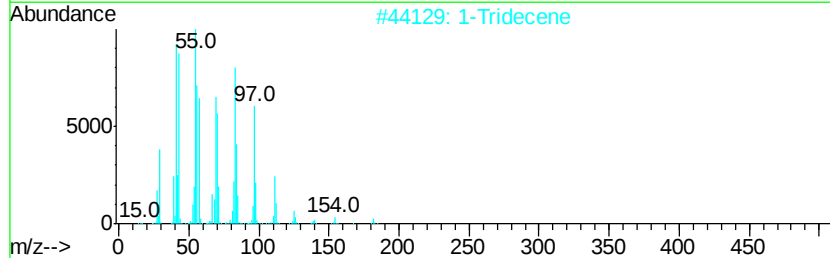
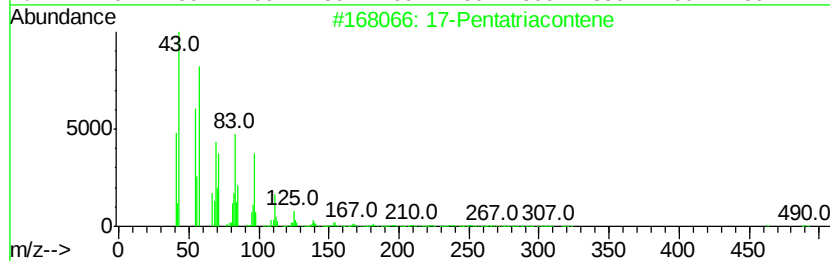
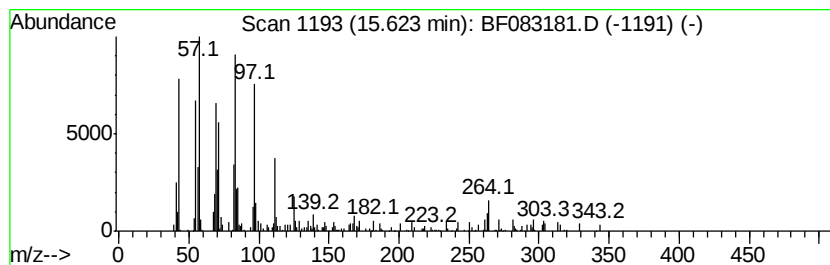
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF112115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 15 17-Pentatriacontene Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.62	3.53 ng	187174	Perylene-d12	15.15	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	17-Pentatriacontene	491	C35H70	006971-40-0	74
2	1-Tridecene	182	C13H26	002437-56-1	64
3	Cvclopropane, nonyl-	168	C12H24	074663-85-7	60
4	Cvclopropane, 1-(1,2-dimethylpro...	252	C18H36	041977-42-8	52
5	Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	49



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF112215\
Data File : BF083181.D
Acq On : 23 Nov 2015 2:04
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ClientSampleId :
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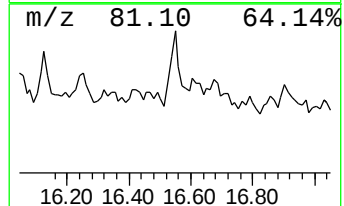
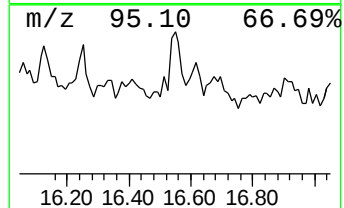
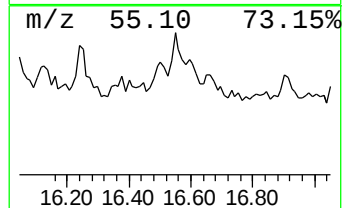
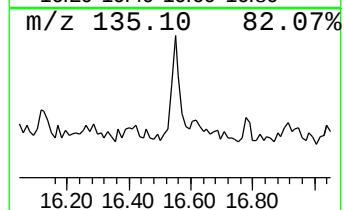
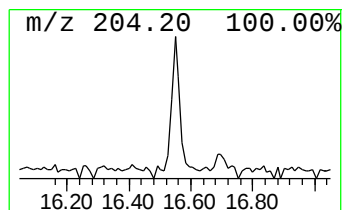
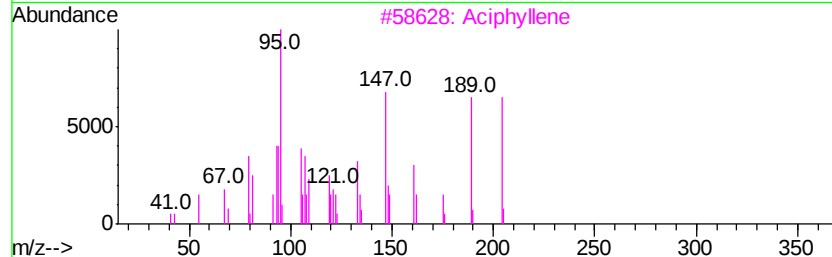
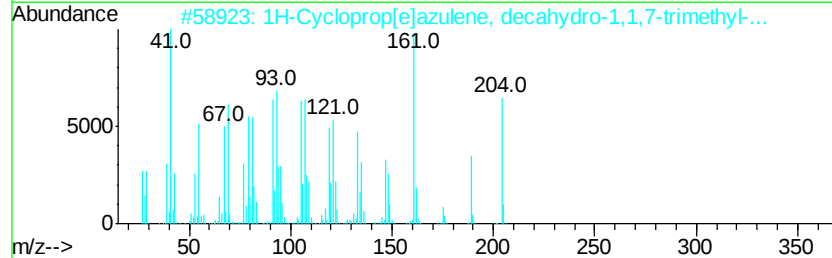
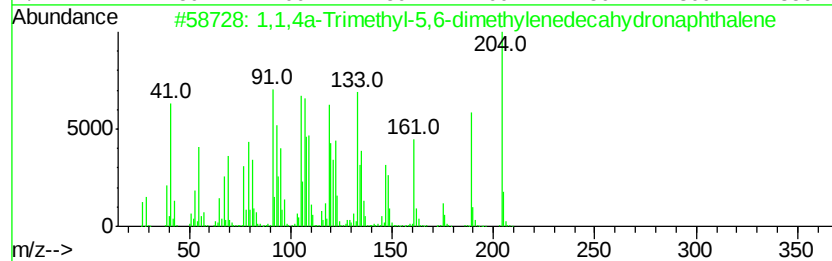
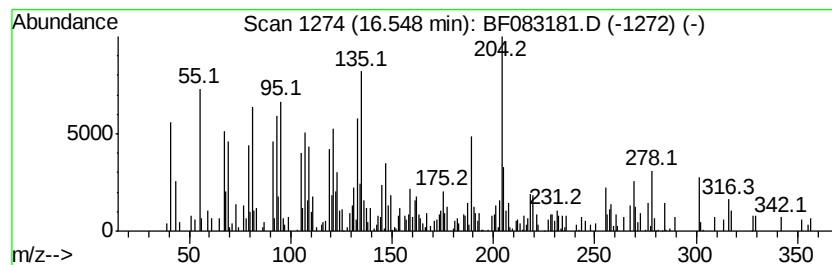
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 16 1,1,4a-Trimethyl-5,6-dimeth... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.55	4.53 ng	240623	Perylene-d12	15.15

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,1,4a-Trimethyl-5,6-dimethylene...	204	C15H24	1000193-60-8	62
2			1H-Cycloprop[elazulene, decahydr...	204	C15H24	000489-39-4	58
3			Aciphyllene	204	C15H24	087745-31-1	55
4			1,2,4,8-Tetramethylbicyclo[6.3.0...	204	C15H24	137235-51-9	55
5			Azulene, 1,2,3,4,5,6,7,8-octahyd...	204	C15H24	003691-12-1	53



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF112215\
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Acq On : 23 Nov 2015 2:04
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Sample : G4512-02
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ALS Vial : 67 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
FILL-DIRT-2

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF112115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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unknown6.42	6.42	81.2	ng	5668180	1	6.65	1395430	20.0
Eicosane	10.14	3.7	ng	333234	3	9.68	1808510	20.0
Pentadecane, 2,6,...	10.35	2.3	ng	203454	3	9.68	1808510	20.0
Tridecanoic acid	11.71	10.2	ng	1002420	4	11.16	1957740	20.0
Tetradecanoic acid	12.49	4.1	ng	431639	5	13.78	2084460	20.0
1-Tetracosanol	13.66	6.9	ng	715810	5	13.78	2084460	20.0
1-Hentetracontanol	14.29	4.1	ng	423047	5	13.78	2084460	20.0
Heneicosane	15.58	5.3	ng	279952	6	15.15	1061590	20.0
17-Pentatriacontene	15.62	3.5	ng	187174	6	15.15	1061590	20.0
1,1,4a-Trimethyl-...	16.55	4.5	ng	240623	6	15.15	1061590	20.0