

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF051019\
 Data File : BF114126.D
 Acq On : 10 May 2019 17:55
 Operator : JU/SJ
 Sample : K2764-08
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 P021-SS010-0002-01

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 3 % of largest Peak

Max Peaks: 100

Peak Location: TOP

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

Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF051019.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	2.152	5	11	19	rVB	1500166	1978923	21.29%	2.543%
2	5.481	572	577	581	rBV	8029695	8806539	94.74%	11.315%
3	6.492	743	749	752	rBV	8528984	9295478	100.00%	11.943%
4	6.628	767	772	775	rBV	8725663	8463085	91.05%	10.873%
5	6.851	806	810	813	rBV	1814556	1395018	15.01%	1.792%
6	7.010	833	837	840	rBV	5809412	4890864	52.62%	6.284%
7	7.410	900	905	908	rBV	4068920	4281865	46.06%	5.501%
8	8.128	1023	1027	1031	rBV	1863316	1744061	18.76%	2.241%
9	9.204	1205	1210	1214	rBV	6372921	6348167	68.29%	8.156%
10	9.598	1272	1277	1280	rBV	880413	765822	8.24%	0.984%
11	9.886	1319	1326	1329	rBV	2026015	1891557	20.35%	2.430%
12	10.675	1455	1460	1463	rBV	4433523	4126869	44.40%	5.302%
13	11.369	1573	1578	1581	rBV	2270102	1924234	20.70%	2.472%
14	11.898	1664	1668	1679	rBV	1221245	1291765	13.90%	1.660%
15	12.580	1780	1784	1786	rBV	175530	168634	1.81%	0.217%
16	12.680	1797	1801	1809	rBV	276979	374205	4.03%	0.481%
17	12.804	1818	1822	1825	rBV	158800	156333	1.68%	0.201%
18	12.951	1842	1847	1850	rBV	6706187	6182223	66.51%	7.943%
19	13.851	1996	2000	2013	rBV	984810	1361135	14.64%	1.749%
20	14.015	2023	2028	2031	rBV	2002810	2098569	22.58%	2.696%
21	14.186	2054	2057	2060	rBV	153306	144083	1.55%	0.185%
22	14.504	2108	2111	2117	rBV2	352705	417938	4.50%	0.537%
23	14.815	2160	2164	2168	rVB	493169	552048	5.94%	0.709%
24	14.963	2186	2189	2192	rBV	131363	117980	1.27%	0.152%
25	15.074	2205	2208	2216	rVB	80535	144311	1.55%	0.185%
26	15.157	2217	2222	2226	rBV	1242669	1548190	16.66%	1.989%
27	15.504	2276	2281	2284	rBV	1692481	2082148	22.40%	2.675%
28	15.539	2284	2287	2294	rVB	1150456	1479261	15.91%	1.901%
29	15.739	2318	2321	2326	rVB3	127145	157974	1.70%	0.203%
30	15.974	2355	2361	2366	rBV	1177556	1778780	19.14%	2.285%
31	16.027	2367	2370	2375	rVV2	100333	174537	1.88%	0.224%
32	16.480	2441	2447	2453	rBV	396272	734364	7.90%	0.944%
33	16.768	2492	2496	2502	rVB4	105362	176492	1.90%	0.227%
34	17.068	2543	2547	2556	rVB	175686	345965	3.72%	0.444%

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

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

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35	17.315	2586	2589	2598	rVB3	153043	306107	3.29%	0.393%
36	17.574	2628	2633	2634	rBV5	81248	127750	1.37%	0.164%

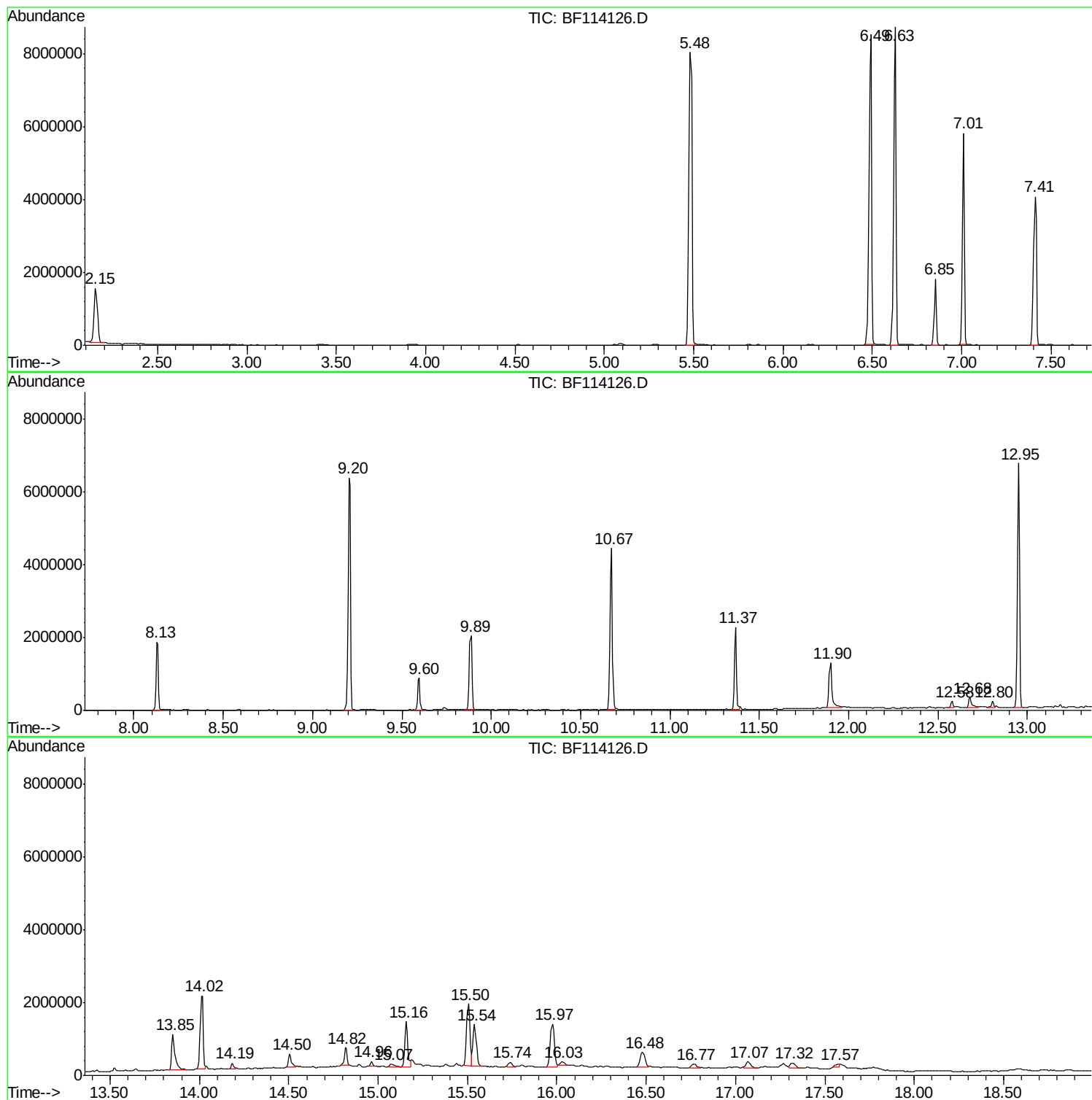
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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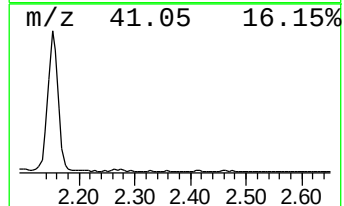
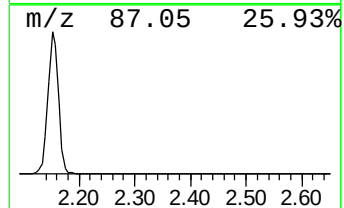
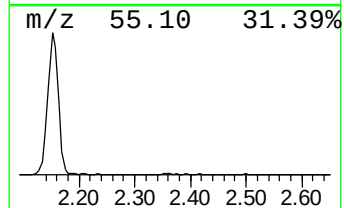
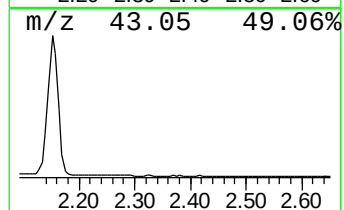
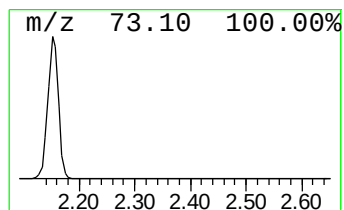
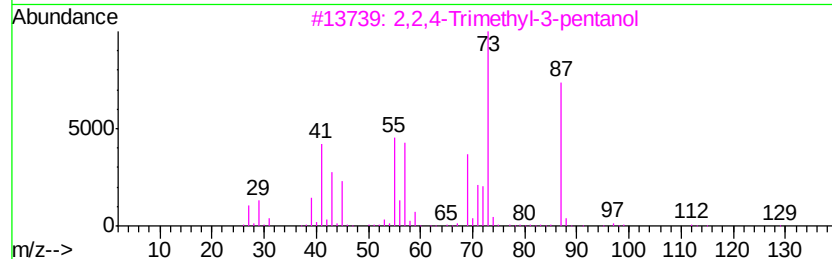
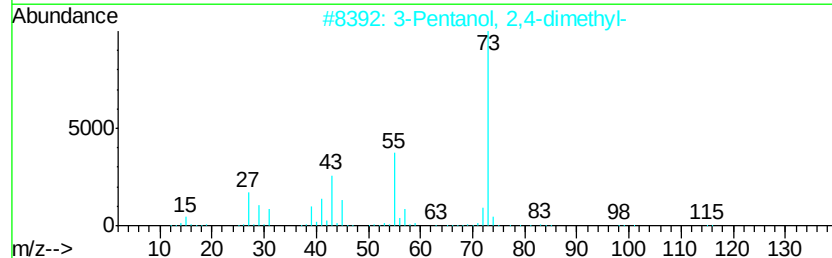
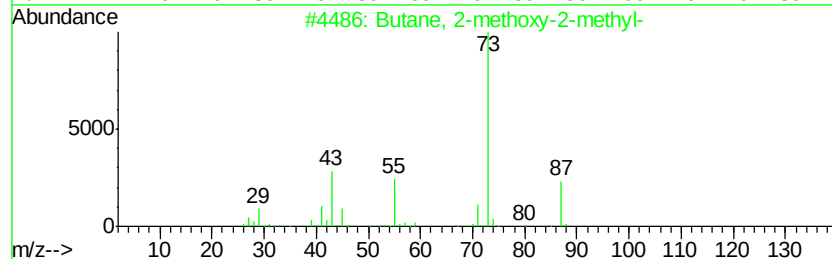
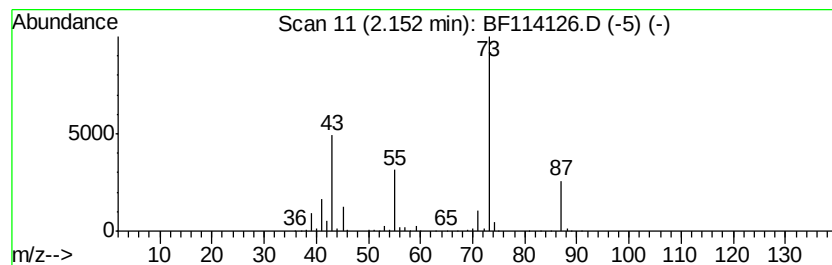
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.15	28.37 ng	1978920	1,4-Dichlorobenzene-d4	6.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2		3-Pentanol, 2,4-dimethyl-	116	C7H16O	000600-36-2	43
3		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
4		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	23
5		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12



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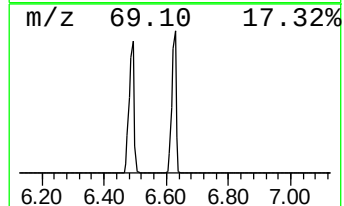
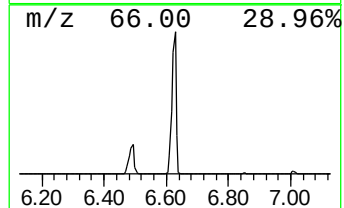
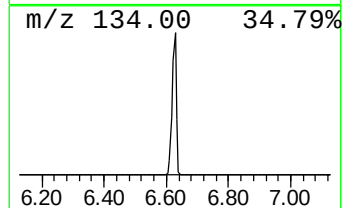
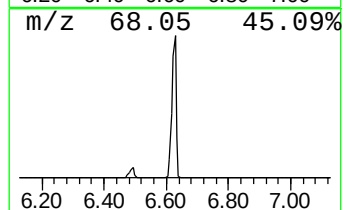
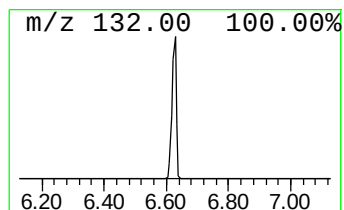
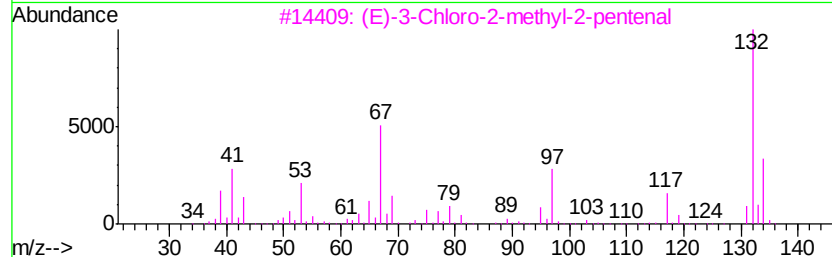
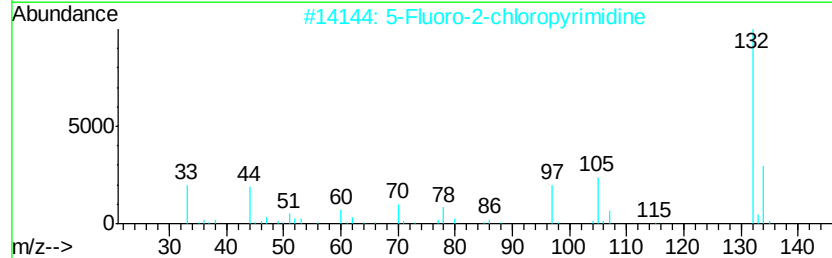
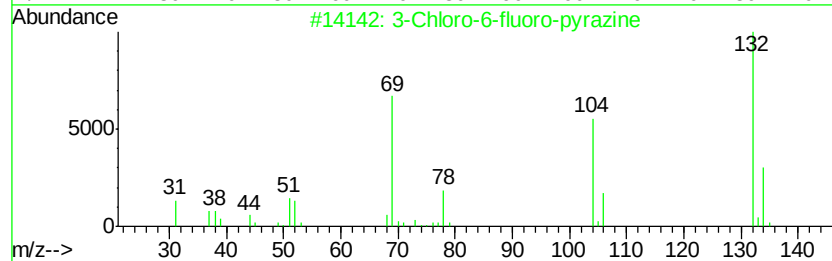
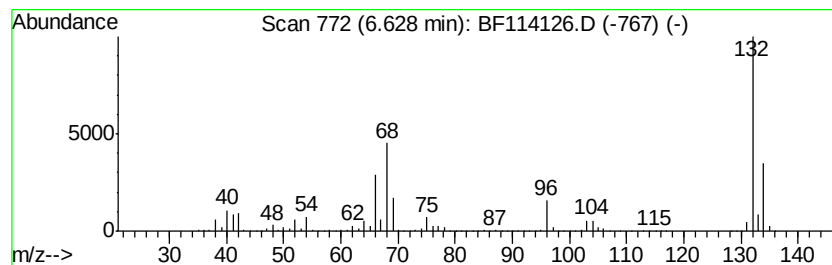
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 2 unknown6.63 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.63	121.33 ng	8463090	1,4-Dichlorobenzene-d4	6.85

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		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
4		Xenon	132	Xe	007440-63-3	9
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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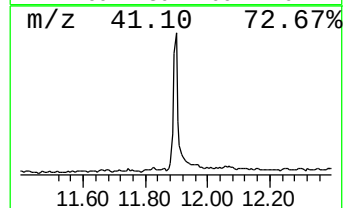
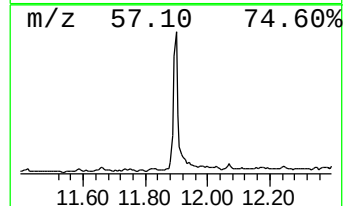
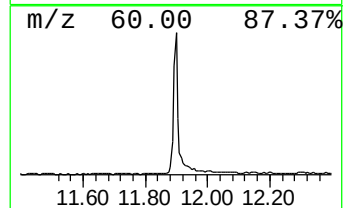
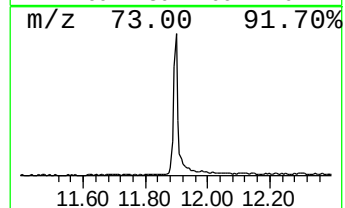
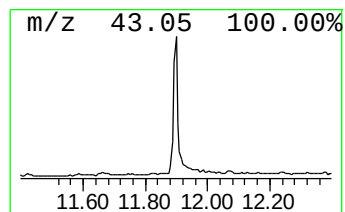
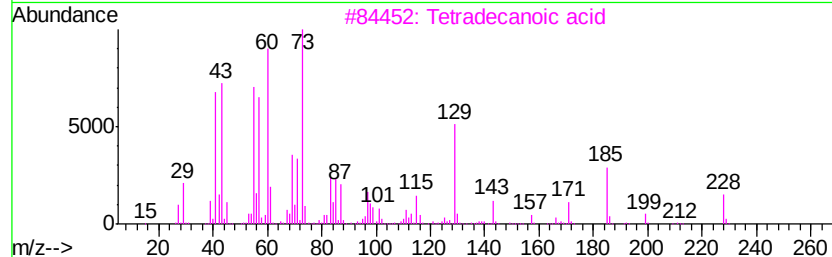
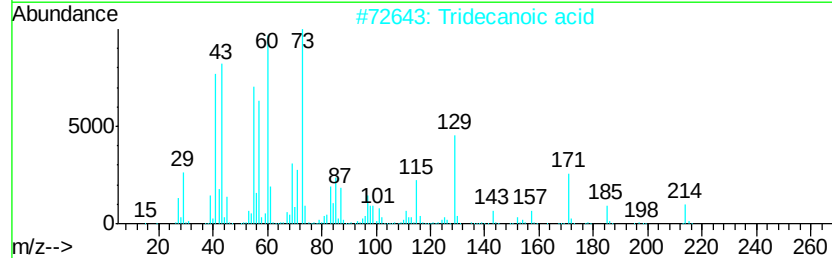
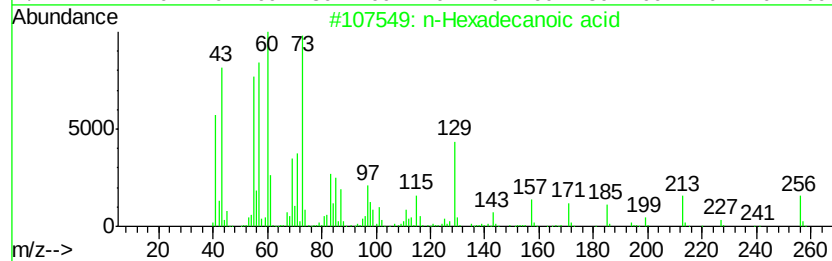
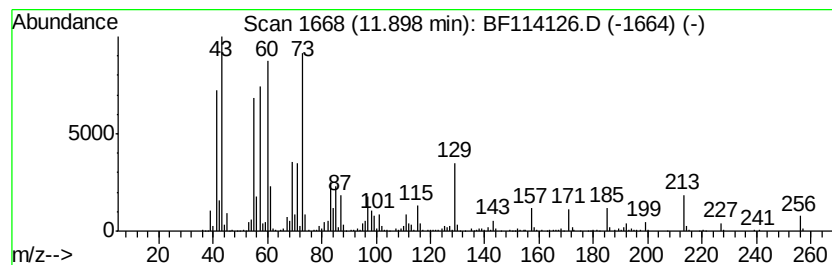
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 4 n-Hexadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.90	13.43 ng	1291770	Phenanthrene-d10	11.37

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		Tridecanoic acid	214	C13H26O2	000638-53-9	93
3		Tetradecanoic acid	228	C14H28O2	000544-63-8	87
4		Pentadecanoic acid	242	C15H30O2	001002-84-2	87
5		Undecanoic acid	186	C11H22O2	000112-37-8	64



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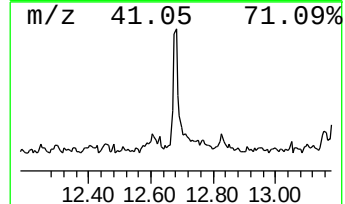
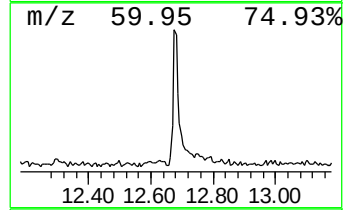
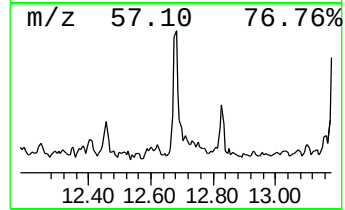
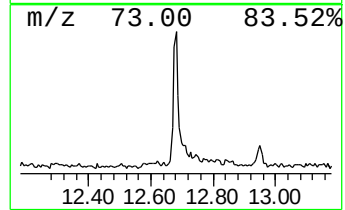
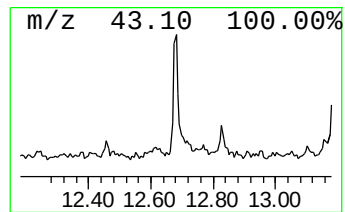
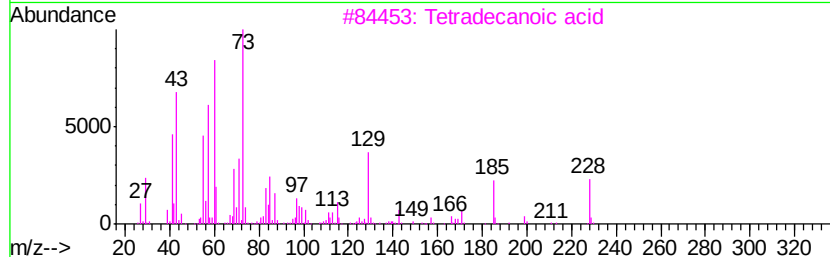
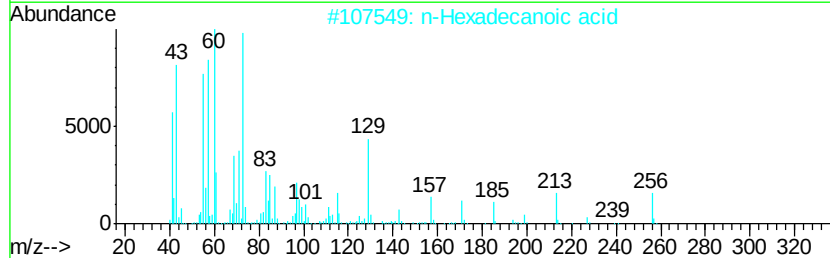
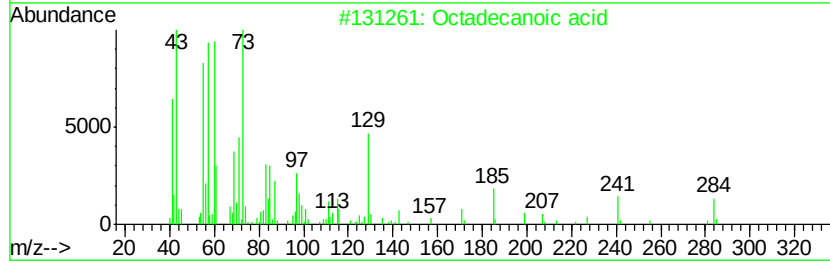
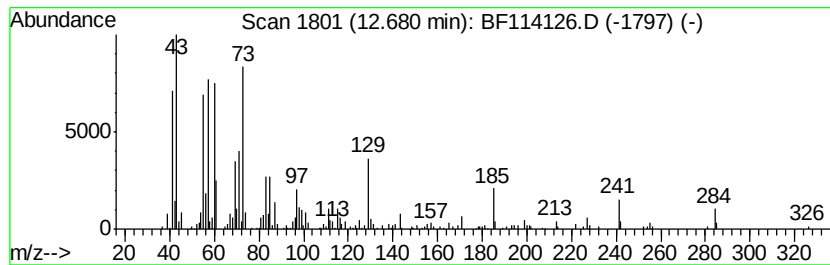
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 5 Octadecanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD		R.T.
12.68	3.89 ng	374205	Phenanthrene-d10		11.37
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	87
3	Tetradecanoic acid	228	C14H28O2	000544-63-8	68
4	Pentadecanoic acid	242	C15H30O2	001002-84-2	68
5	Tridecanoic acid	214	C13H26O2	000638-53-9	62



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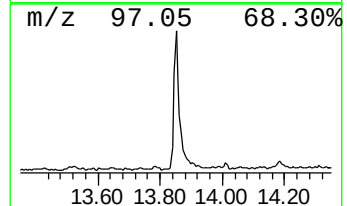
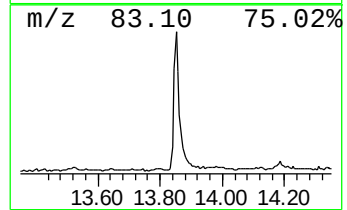
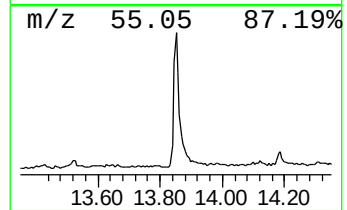
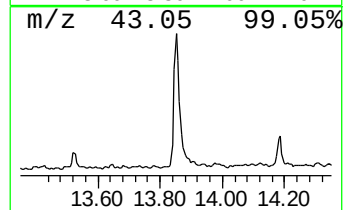
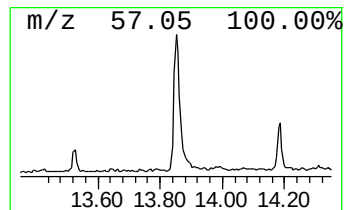
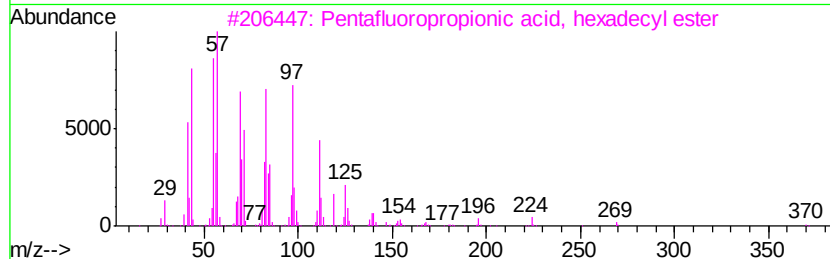
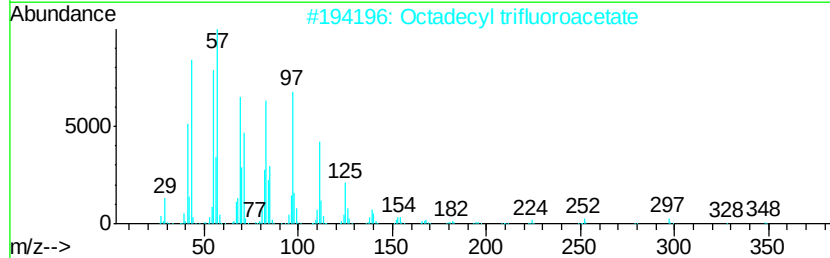
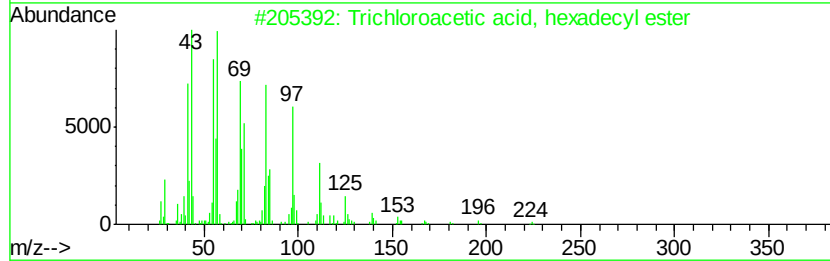
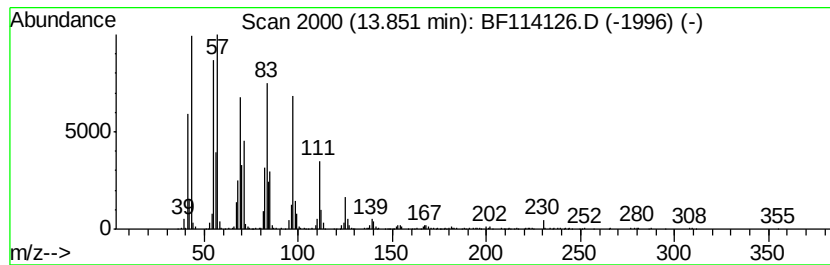
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Peak Number 6 Trichloroacetic acid, hexad... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.		
13.85	12.97 ng	1361140	Chrysene-d12	14.02		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	94
2		Octadecyl trifluoroacetate	366	C20H37F3O2	079392-43-1	94
3		Pentafluoropropionic acid, hexad...	388	C19H33F5O2	006222-07-7	94
4		Pentafluoropropionic acid, octad...	416	C21H37F5O2	959261-25-7	94
5		Behenic alcohol	326	C22H46O	000661-19-8	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF051019\
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Acq On : 10 May 2019 17:55
Operator : JU/SJ
Sample : K2764-08
Misc :
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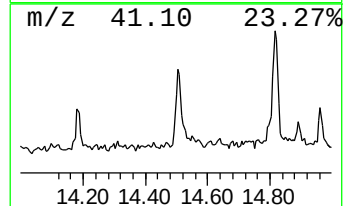
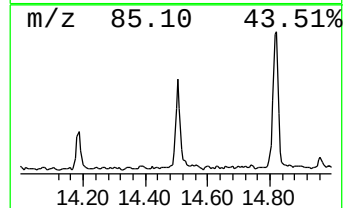
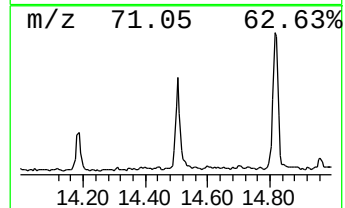
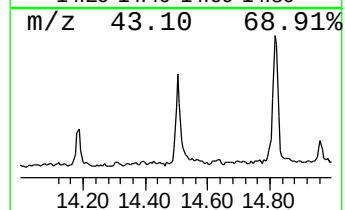
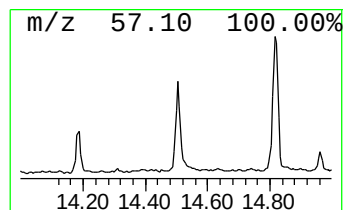
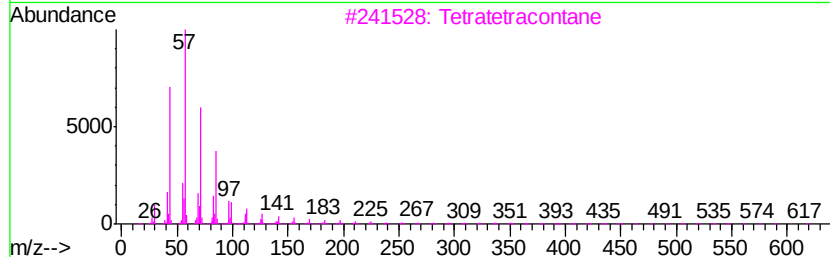
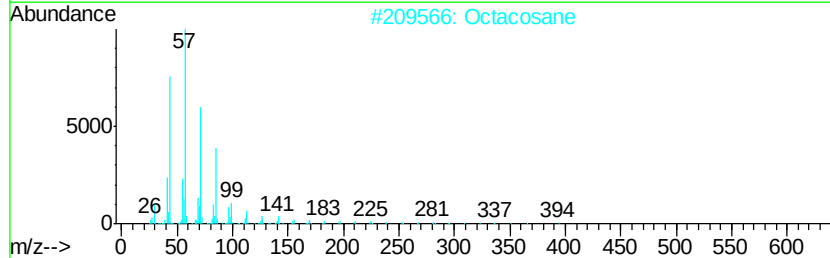
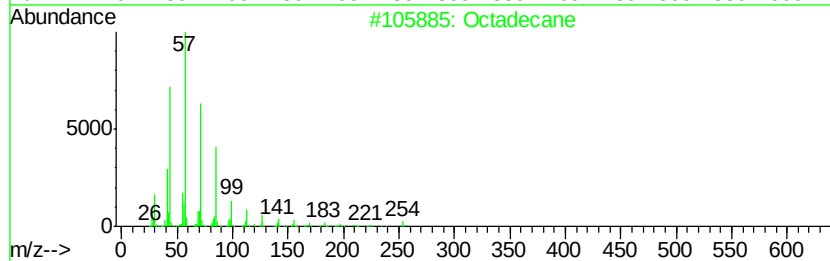
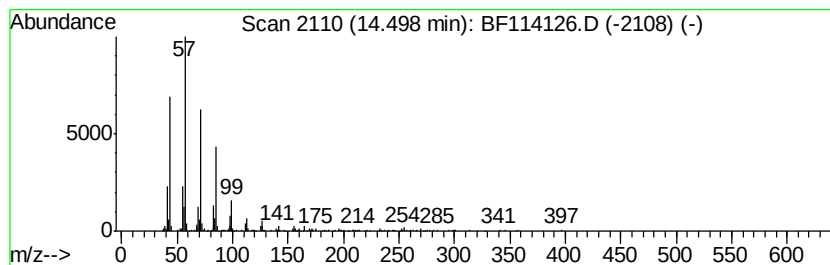
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Octadecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.50	3.98 ng	417938	Chrysene-d12	14.02
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Octadecane	254 C18H38	000593-45-3	93
2	Octacosane	394 C28H58	000630-02-4	91
3	Tetratetracontane	619 C44H90	007098-22-8	91
4	Pentacosane	352 C25H52	000629-99-2	91
5	Hexatriacontane	507 C36H74	000630-06-8	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF051019\
Data File : BF114126.D
Acq On : 10 May 2019 17:55
Operator : JU/SJ
Sample : K2764-08
Misc :
ALS Vial : 4 Sample Multiplier: 1

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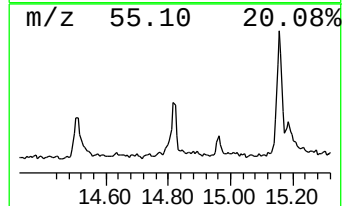
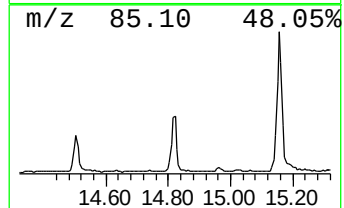
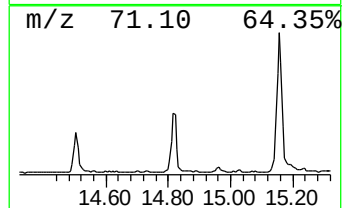
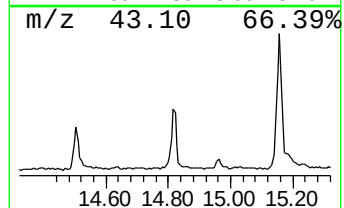
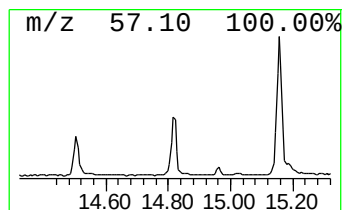
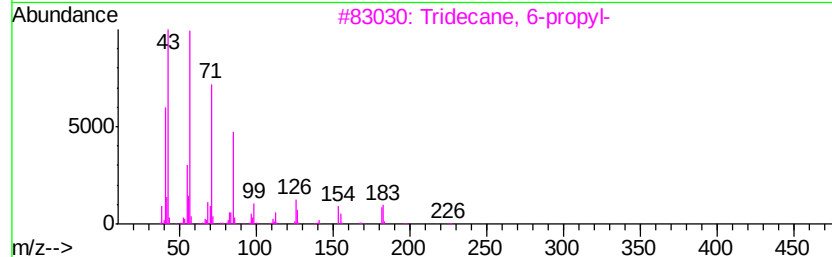
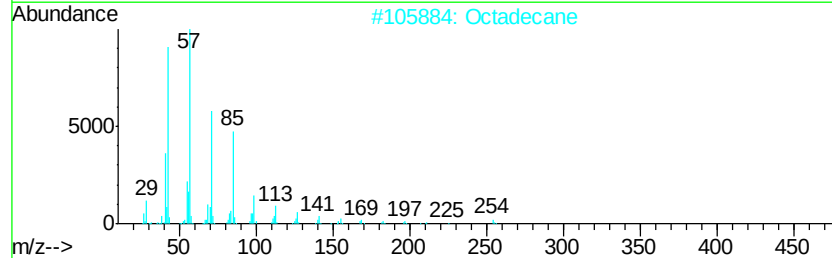
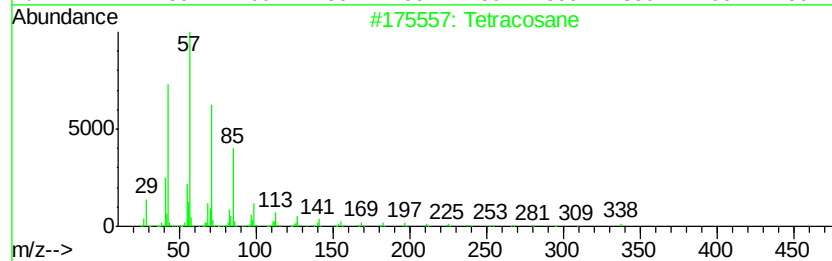
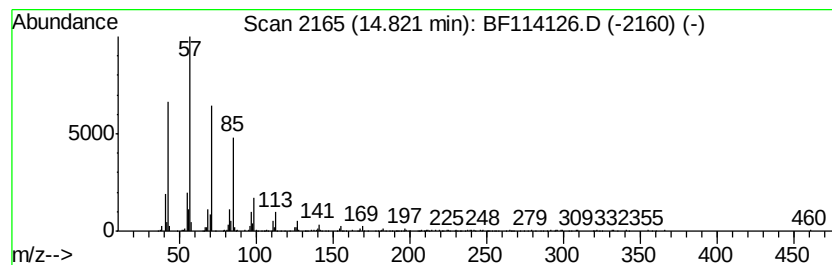
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Tetracosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.82	5.30 ng	552048	Perylene-d12	15.50

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetracosane	338	C24H50	000646-31-1	94
2		Octadecane	254	C18H38	000593-45-3	91
3		Tridecane, 6-propyl-	226	C16H34	055045-10-8	90
4		Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	90
5		Heptacosane	380	C27H56	000593-49-7	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF051019\
 Data File : BF114126.D
 Acq On : 10 May 2019 17:55
 Operator : JU/SJ
 Sample : K2764-08
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Instrument :
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 ClientSampleId :
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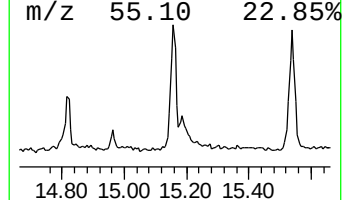
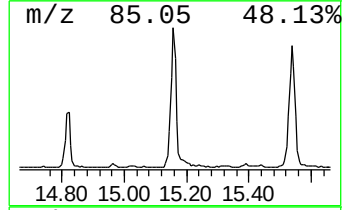
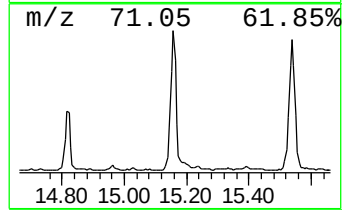
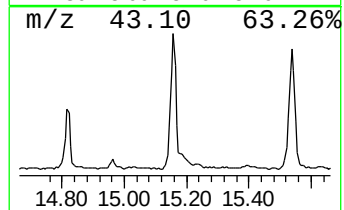
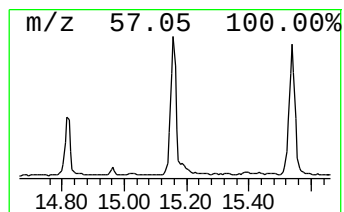
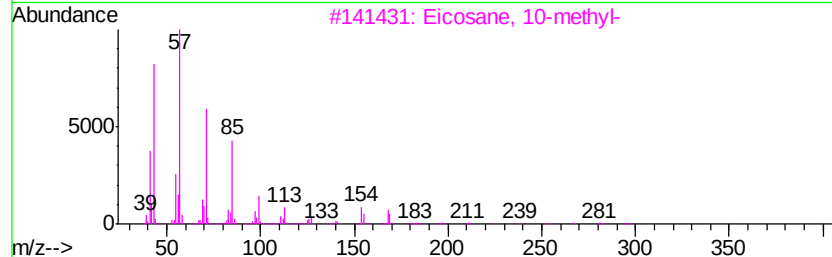
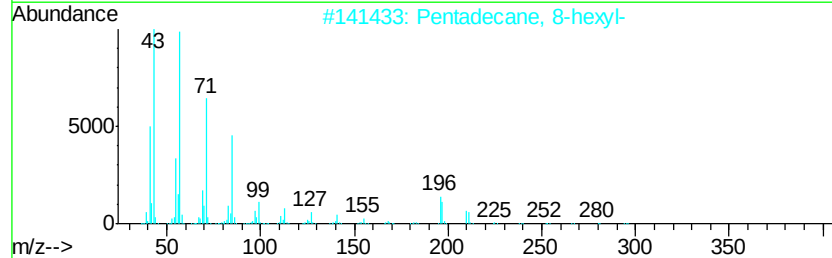
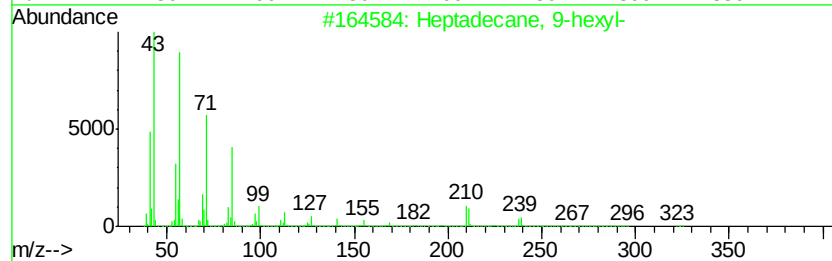
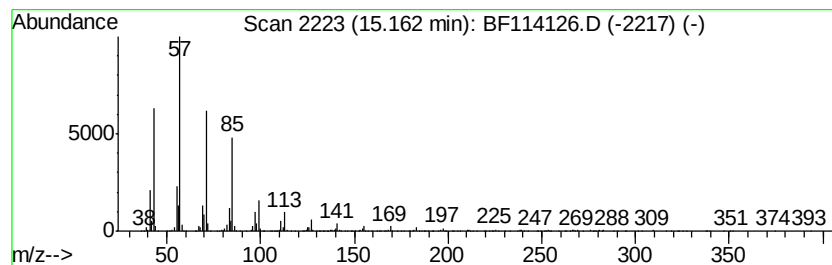
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 9 Heptadecane, 9-hexyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.16	14.87 ng	1548190	Perylene-d12	15.50

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecane, 9-hexyl-	324	C23H48	055124-79-3	94
2		Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	94
3		Eicosane, 10-methyl-	296	C21H44	054833-23-7	93
4		Hexacosane	366	C26H54	000630-01-3	91
5		Pentacosane	352	C25H52	000629-99-2	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF051019\
Data File : BF114126.D
Acq On : 10 May 2019 17:55
Operator : JU/SJ
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Misc :
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Instrument :
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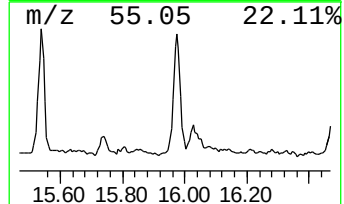
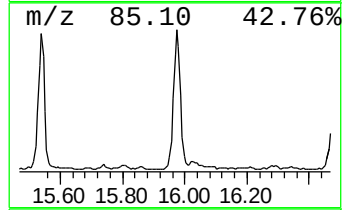
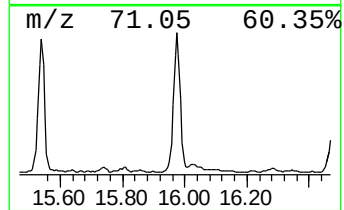
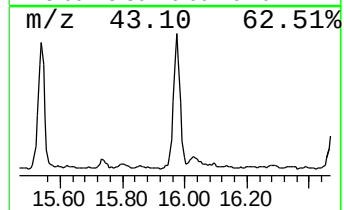
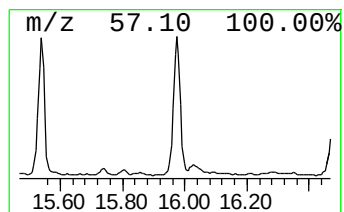
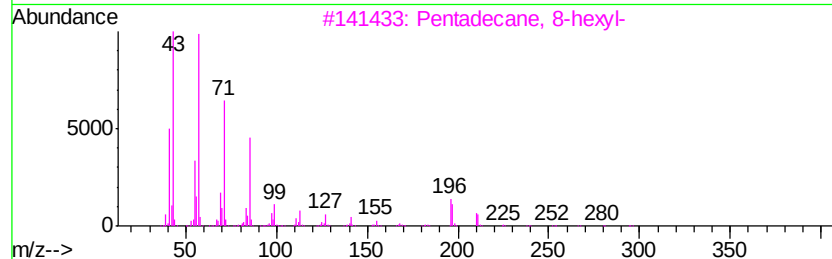
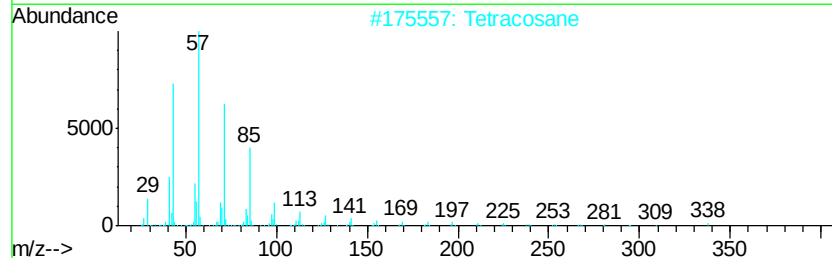
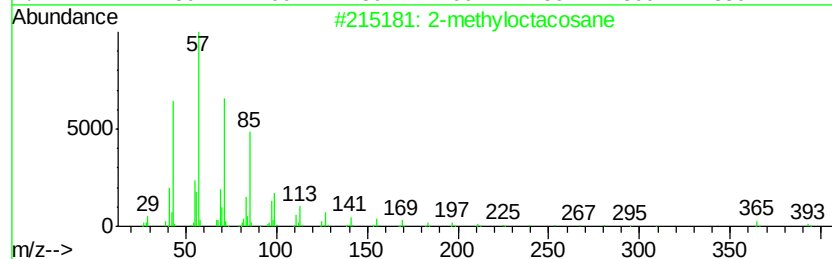
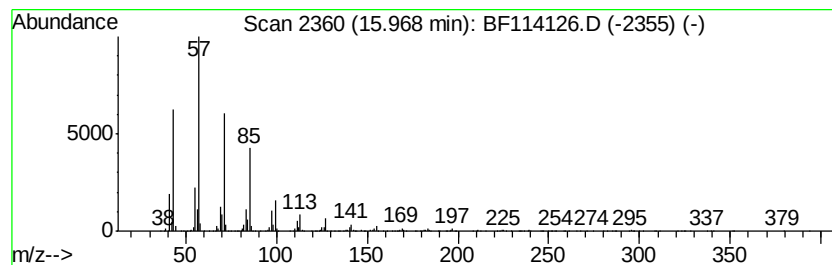
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 2-methyloctacosane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.97	17.09 ng	1778780	Perylene-d12	15.50

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-methyloctacosane	408	C29H60	1000376-72-8	91
2		Tetracosane	338	C24H50	000646-31-1	91
3		Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	91
4		Eicosane, 7-hexyl-	366	C26H54	055333-99-8	91
5		Eicosane, 10-methyl-	296	C21H44	054833-23-7	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF051019\
Data File : BF114126.D
Acq On : 10 May 2019 17:55
Operator : JU/SJ
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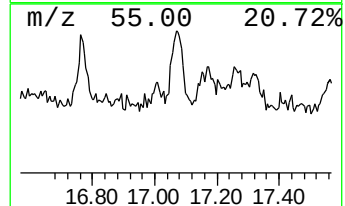
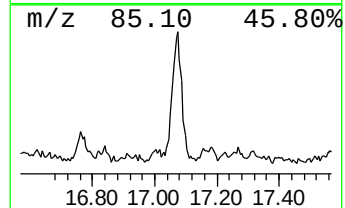
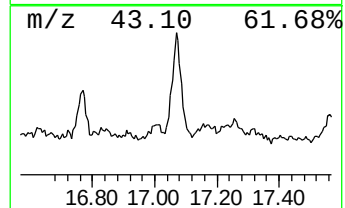
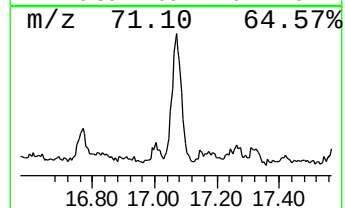
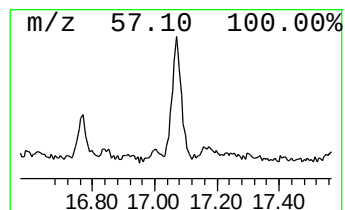
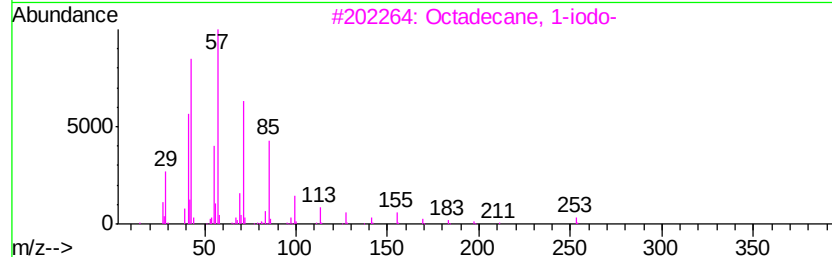
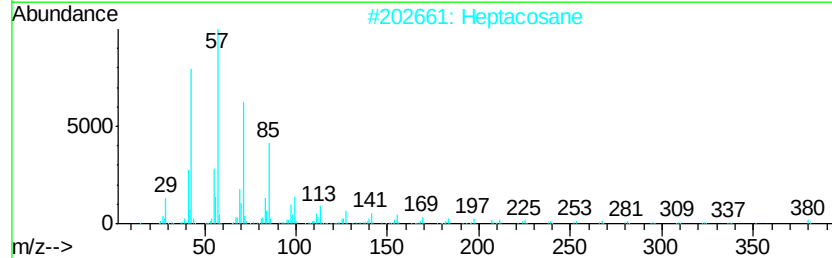
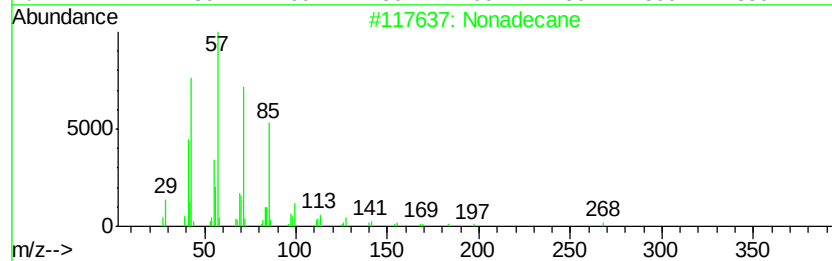
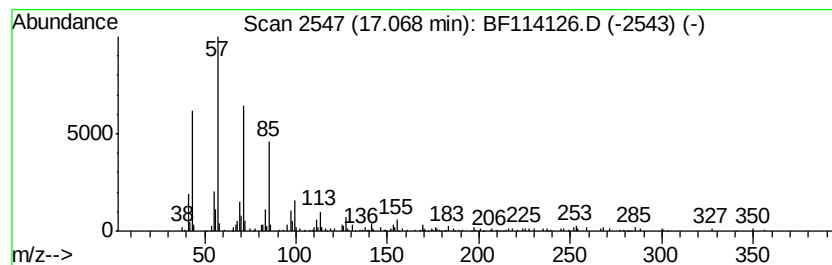
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Nonadecane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.07	3.32 ng	345965	Perylene-d12	15.50

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonadecane	268	C19H40	000629-92-5	93
2		Heptacosane	380	C27H56	000593-49-7	90
3		Octadecane, 1-iodo-	380	C18H37I	000629-93-6	87
4		Heneicosane	296	C21H44	000629-94-7	87
5		Hexacosane	366	C26H54	000630-01-3	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF051019\
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Acq On : 10 May 2019 17:55
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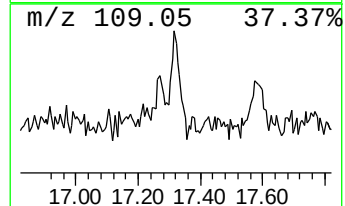
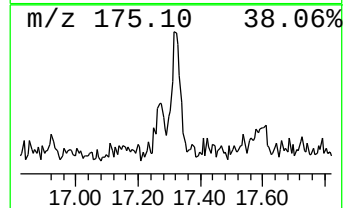
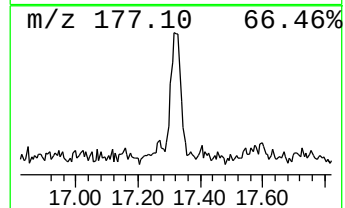
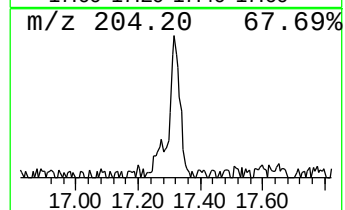
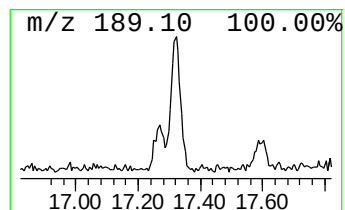
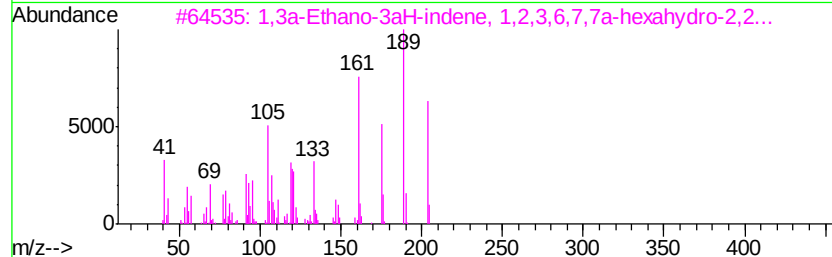
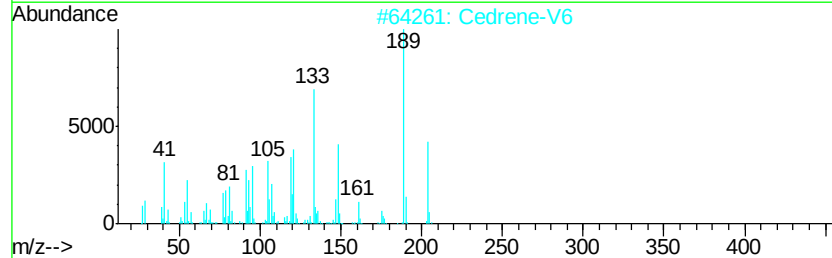
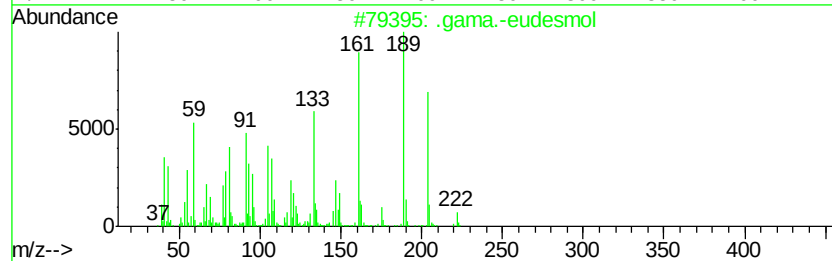
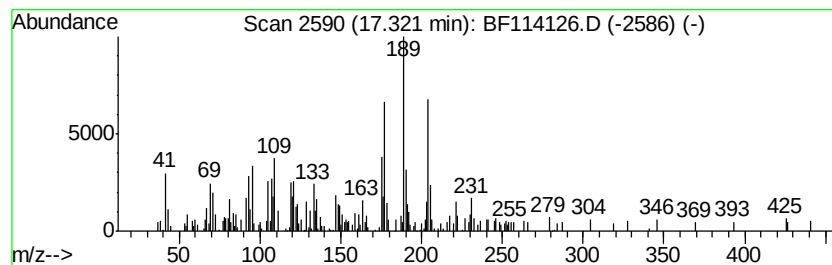
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 .gama.-eudesmol Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.32	2.94 ng	306107	Perylene-d12	15.50

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	.gama.-eudesmol	222	C15H26O	1000374-18-5	64
2		Cedrene-V6	204	C15H24	1000162-76-8	55
3		1,3a-Ethano-3aH-indene, 1,2,3,6,...	204	C15H24	004545-68-0	52
4		1H-Cyclopropa[a]naphthalene, 1a,...	204	C15H24	000489-29-2	46
5		.delta.-Selinene	204	C15H24	028624-23-9	43



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF051019\
Data File : BF114126.D
Acq On : 10 May 2019 17:55
Operator : JU/SJ
Sample : K2764-08
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF051019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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unknown6.63	6.63	121.3	ng	8463090	1	6.85	1395020	20.0
n-Hexadecanoic acid	11.90	13.4	ng	1291770	4	11.37	1924230	20.0
Octadecanoic acid	12.68	3.9	ng	374205	4	11.37	1924230	20.0
Trichloroacetic a...	13.85	13.0	ng	1361140	5	14.02	2098570	20.0
Octadecane	14.50	4.0	ng	417938	5	14.02	2098570	20.0
Tetracosane	14.82	5.3	ng	552048	6	15.50	2082150	20.0
Heptadecane, 9-he...	15.16	14.9	ng	1548190	6	15.50	2082150	20.0
2-methyloctacosane	15.97	17.1	ng	1778780	6	15.50	2082150	20.0
Nonadecane	17.07	3.3	ng	345965	6	15.50	2082150	20.0
.gama.-eudesmol	17.32	2.9	ng	306107	6	15.50	2082150	20.0