

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG041515\
 Data File : BG016439.D
 Acq On : 16 Apr 2015 2:33
 Operator : TP/IZ
 Sample : G1829-12
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 TE-PMW-1A-11

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_G\METHODS\8270-BG040615.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	2.758	7	12	21	rBV	279874	426218	8.95%	1.550%
2	2.835	21	25	39	rVB	347003	450906	9.47%	1.640%
3	3.076	61	66	75	rBV	136801	176401	3.71%	0.642%
4	4.780	351	356	363	rBV	67390	93321	1.96%	0.339%
5	5.255	431	437	447	rBV	639379	918301	19.29%	3.340%
6	6.854	701	709	723	rBV	434923	697277	14.65%	2.536%
7	7.200	761	768	783	rBV	1264302	1989797	41.80%	7.237%
8	7.664	841	847	855	rBV	298791	450803	9.47%	1.640%
9	7.982	894	901	911	rVB	1053215	1701677	35.75%	6.189%
10	8.822	1037	1044	1052	rBV	939248	1500448	31.52%	5.457%
11	10.449	1313	1321	1328	rBV	466670	758350	15.93%	2.758%
12	12.929	1736	1743	1756	rVB	2625783	3835944	80.59%	13.951%
13	14.310	1970	1978	1986	rBV2	737542	1111912	23.36%	4.044%
14	15.673	2204	2210	2215	rVB	63469	85250	1.79%	0.310%
15	15.802	2225	2232	2244	rBV	1702346	2379855	50.00%	8.655%
16	16.078	2274	2279	2285	rVB	41988	67346	1.41%	0.245%
17	17.048	2438	2444	2451	rBV2	880660	1288893	27.08%	4.688%
18	17.947	2592	2597	2606	rBV	172740	278565	5.85%	1.013%
19	18.035	2608	2612	2622	rVB2	47550	86327	1.81%	0.314%
20	19.086	2787	2791	2793	rBV4	49900	69261	1.46%	0.252%
21	19.233	2812	2816	2825	rVB2	104320	144595	3.04%	0.526%
22	19.498	2857	2861	2866	rVB	60206	70957	1.49%	0.258%
23	19.680	2886	2892	2902	rBV2	3722078	4759772	100.00%	17.311%
24	20.350	3003	3006	3008	rBV3	49893	61672	1.30%	0.224%
25	20.485	3024	3029	3033	rBV2	79501	110751	2.33%	0.403%
26	20.908	3098	3101	3104	rBV	118201	133971	2.81%	0.487%
27	20.943	3104	3107	3118	rVB	372782	502726	10.56%	1.828%
28	21.143	3138	3141	3146	rBV	74674	92760	1.95%	0.337%
29	21.231	3151	3156	3161	rVV	962965	1234038	25.93%	4.488%
30	21.378	3178	3181	3185	rVB	113551	125456	2.64%	0.456%
31	21.760	3242	3246	3250	rVB4	82869	137286	2.88%	0.499%
32	21.807	3251	3254	3266	rVB	105517	135939	2.86%	0.494%
33	22.265	3329	3332	3340	rVB	119628	162486	3.41%	0.591%
34	22.394	3351	3354	3359	rVB	64858	78723	1.65%	0.286%

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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 >
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35	22.770	3415	3418	3422	rVB	69309	93354	1.96%	0.340%
36	23.334	3511	3514	3523	rVB2	60604	107493	2.26%	0.391%
37	23.458	3529	3535	3542	rBV	624322	1176566	24.72%	4.279%

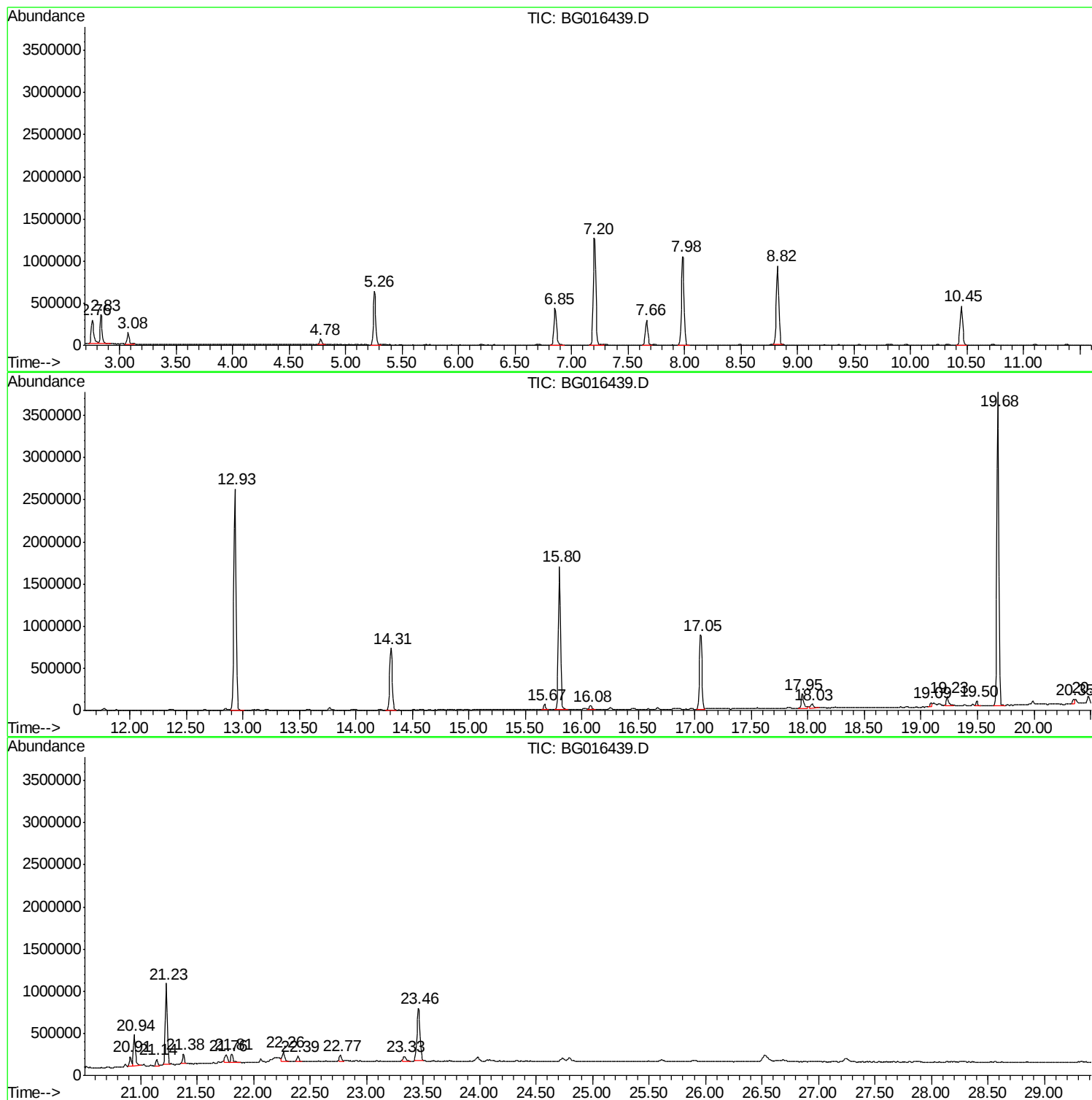
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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG040615.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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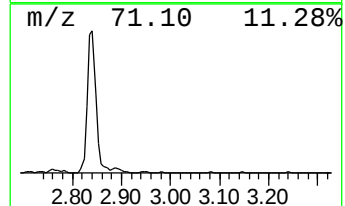
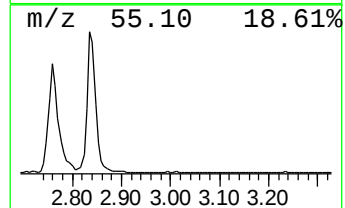
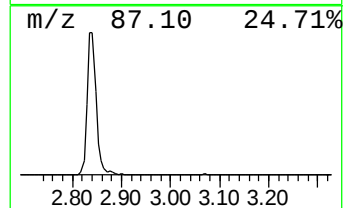
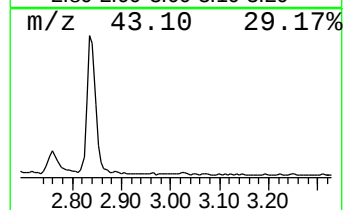
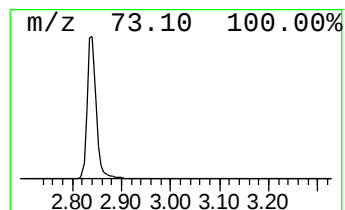
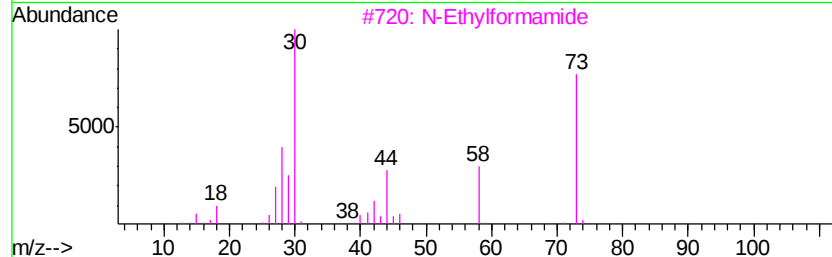
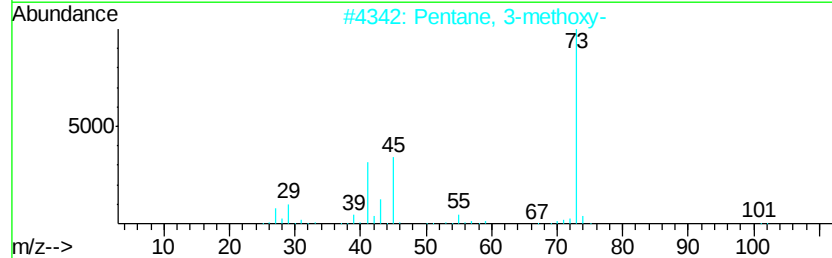
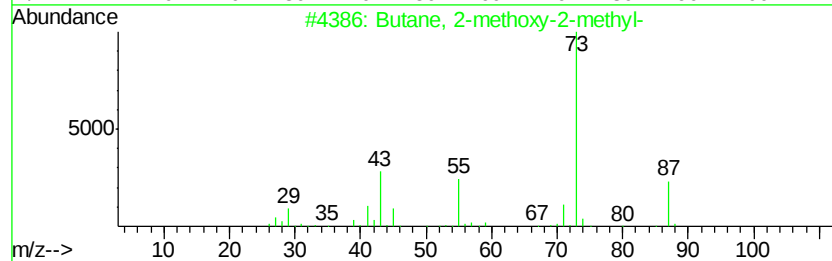
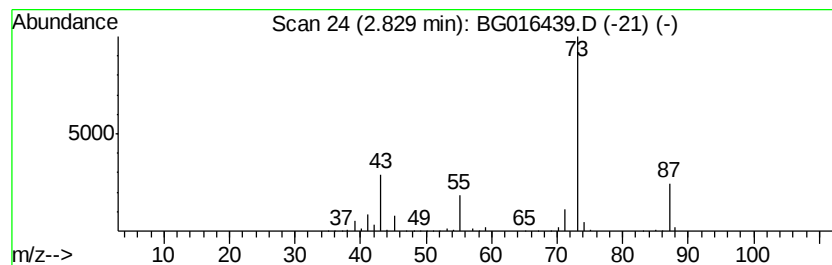
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG040615.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.83	20.00 ng	450906	1,4-Dichlorobenzene-d4	7.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
3		N-Ethylformamide	73	C3H7NO	000627-45-2	9
4		Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9
5		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	7



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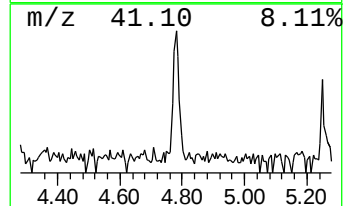
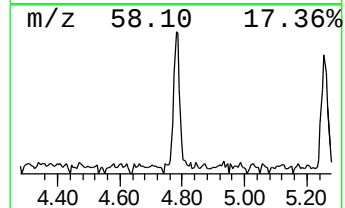
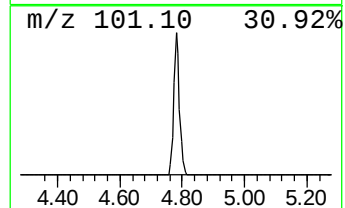
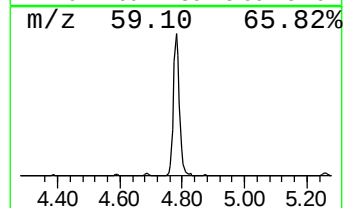
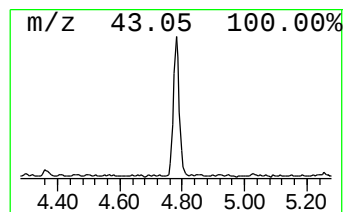
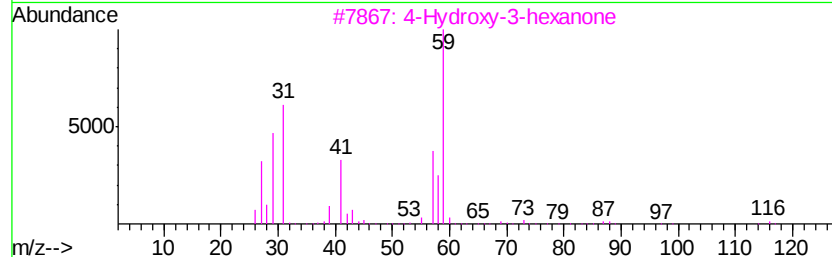
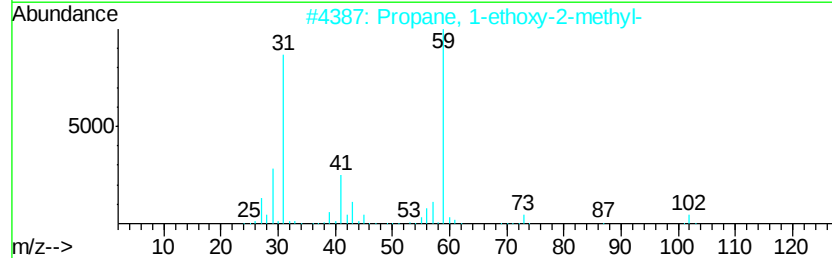
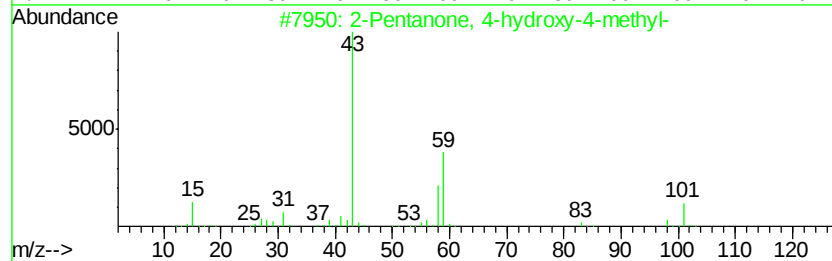
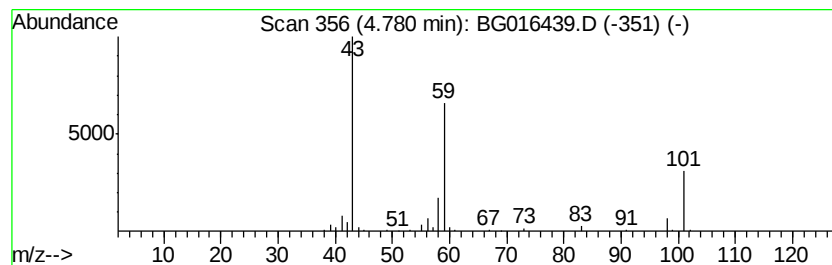
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.78	4.14 ng	93321	1,4-Dichlorobenzene-d4	7.66

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2	Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	25
3	4-Hydroxy-3-hexanone	116	C6H12O2	004984-85-4	23
4	4-Methyl-3-hexanol acetate	158	C9H18O2	084612-71-5	17
5	Pentanamide	101	C5H11NO	000626-97-1	12



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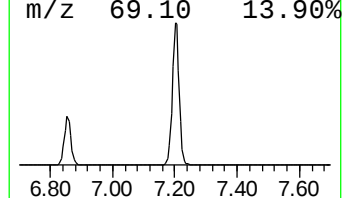
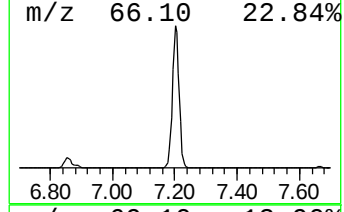
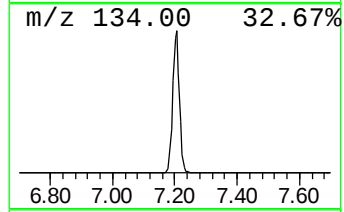
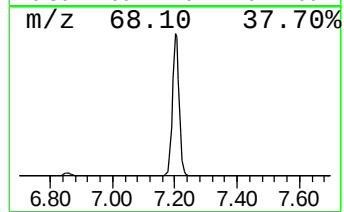
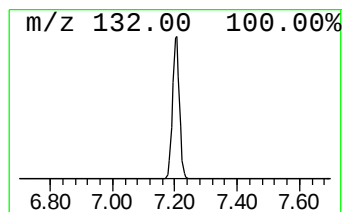
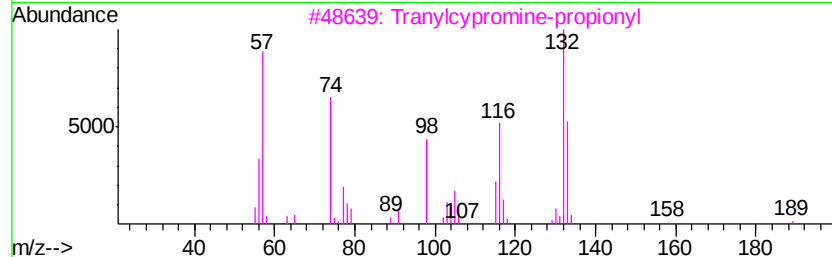
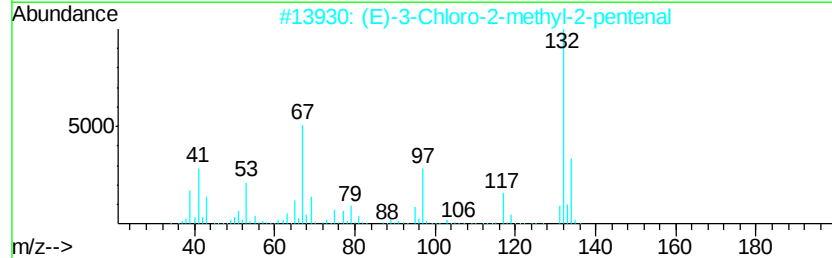
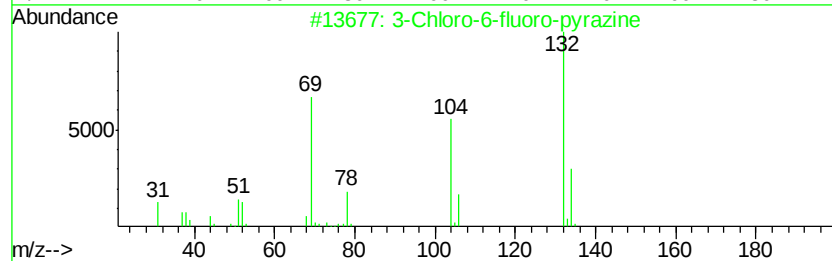
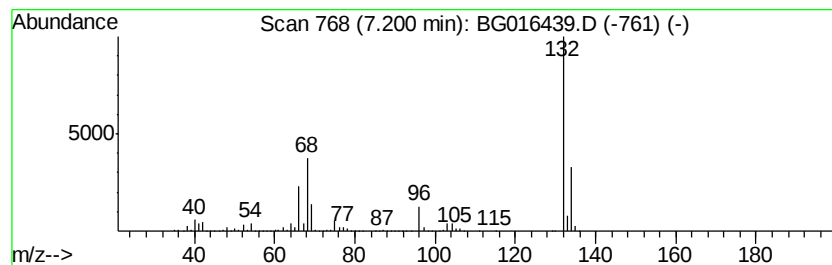
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 unknown7.20 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.20	88.28 ng	1989800	1,4-Dichlorobenzene-d4	7.66

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	38
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	35
3		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
4		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9
5		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9



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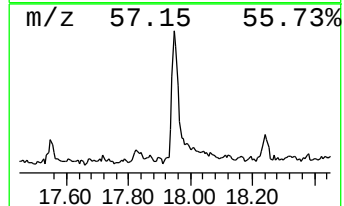
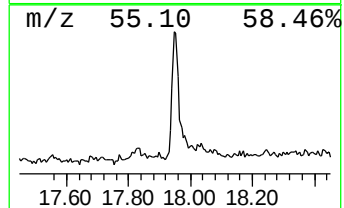
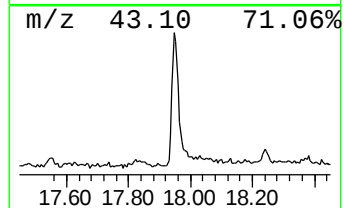
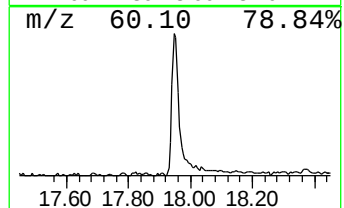
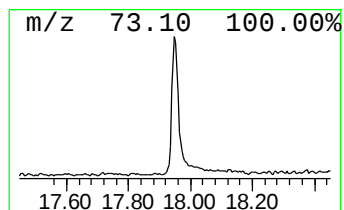
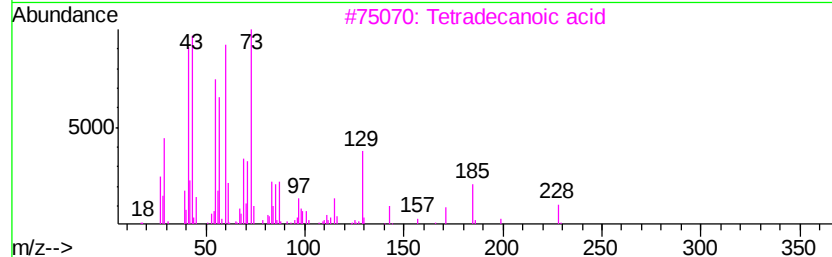
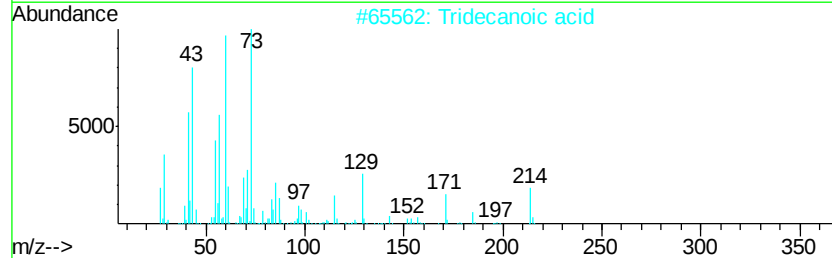
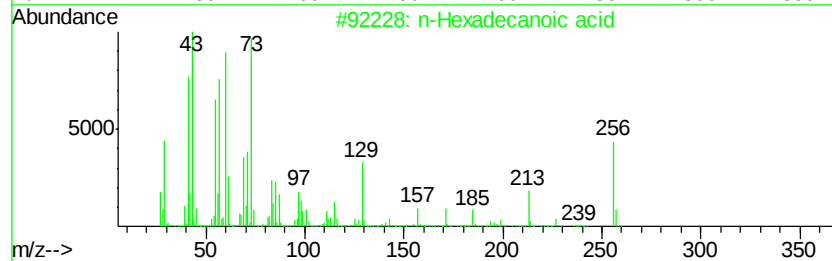
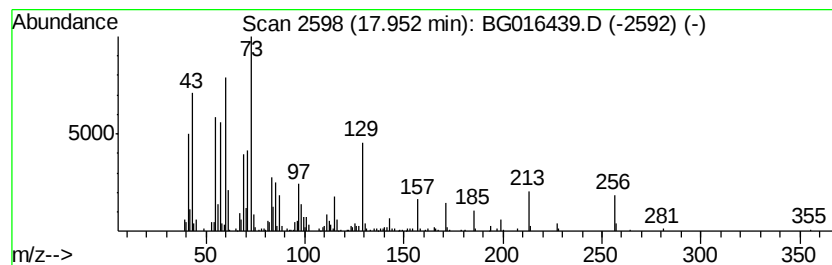
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 n-Hexadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.95	4.32 ng	278565	Phenanthrene-d10	17.05

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2		Tridecanoic acid	214	C13H26O2	000638-53-9	86
3		Tetradecanoic acid	228	C14H28O2	000544-63-8	78
4		Pentadecanoic acid	242	C15H30O2	001002-84-2	78
5		n-Decanoic acid	172	C10H20O2	000334-48-5	58



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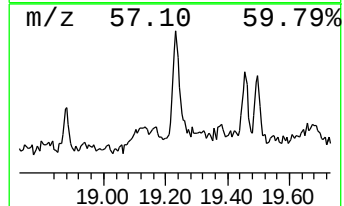
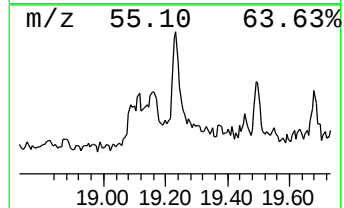
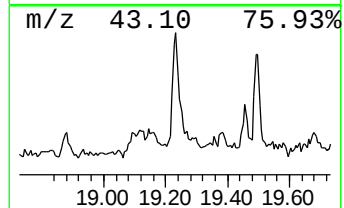
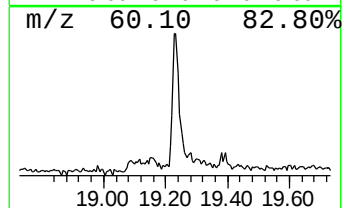
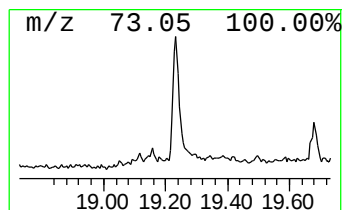
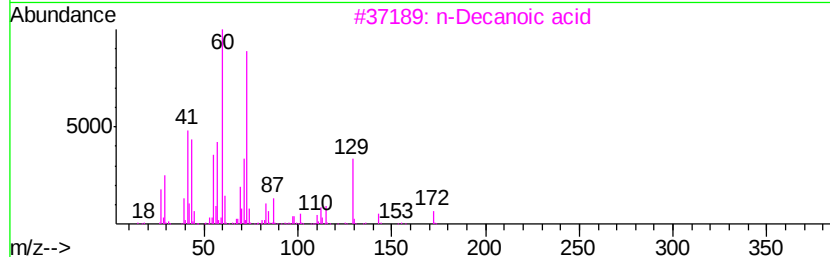
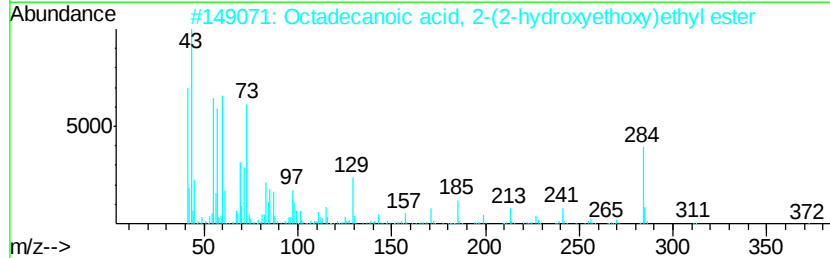
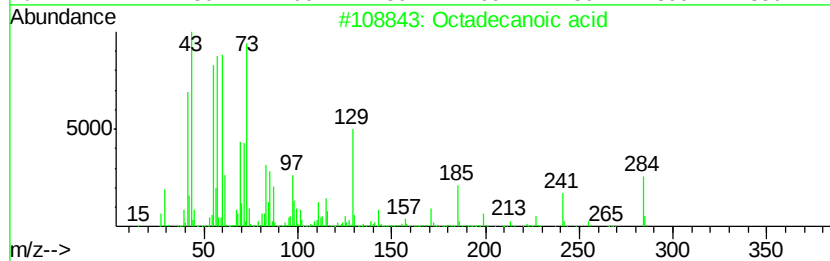
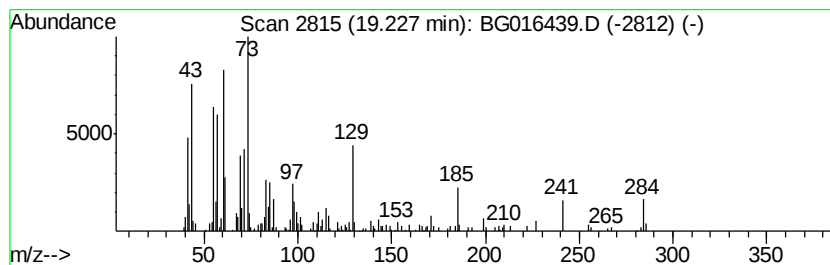
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 8 Octadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.23	2.34 ng	144595	Chrysene-d12	21.23

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2		Octadecanoic acid, 2-(2-hydroxyethyl) ester	372	C22H44O4	000106-11-6	90
3		n-Decanoic acid	172	C10H20O2	000334-48-5	83
4		Pentadecanoic acid	242	C15H30O2	001002-84-2	68
5		Tetradecanoic acid	228	C14H28O2	000544-63-8	53



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG041515\
Data File : BG016439.D
Acq On : 16 Apr 2015 2:33
Operator : TP/IZ
Sample : G1829-12
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
TE-PMW-1A-11

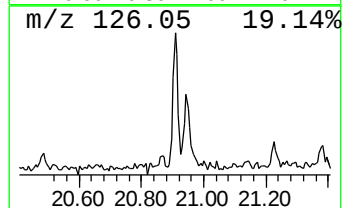
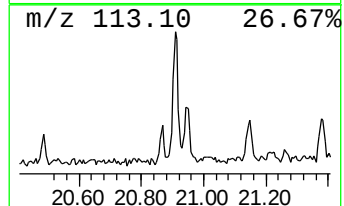
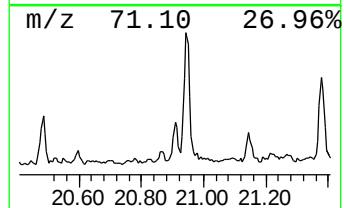
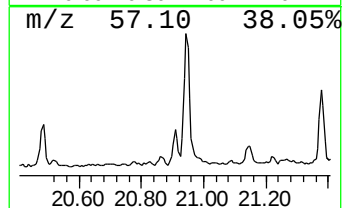
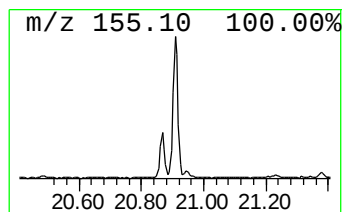
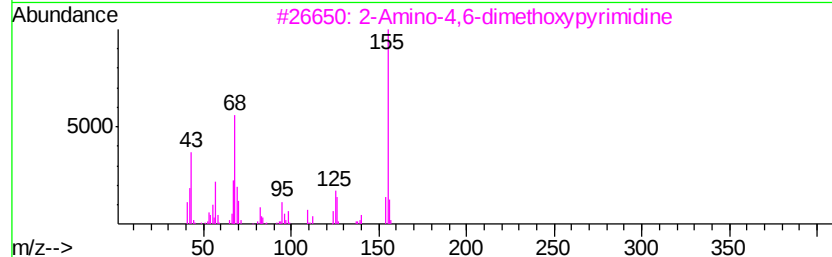
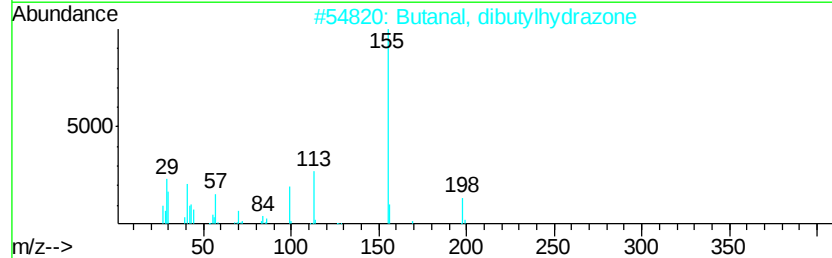
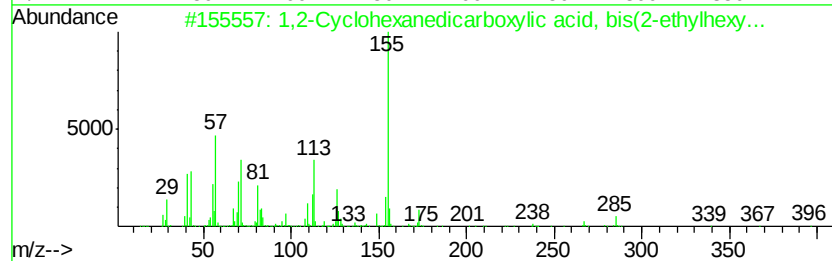
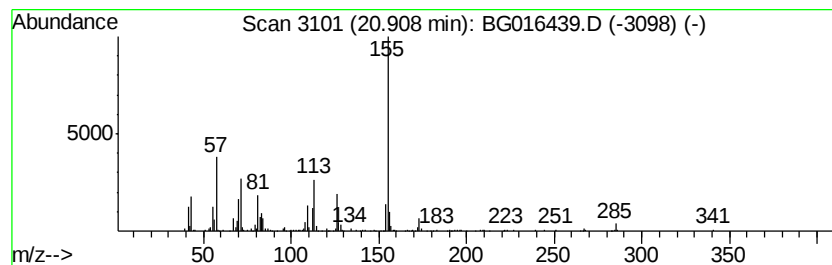
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 1,2-Cyclohexanedicarboxylic... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.91	2.17 ng	133971	Chrysene-d12	21.23

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,2-Cyclohexanedicarboxylic acid...	396	C24H44O4	000084-71-9	62
2		Butanal, dibutylhydrazone	198	C12H26N2	067660-51-9	37
3		2-Amino-4,6-dimethoxypyrimidine	155	C6H9N3O2	036315-01-2	33
4		3,3'-Dimenthol	310	C20H38O2	148552-94-7	28
5		1,4-Dioxaspiro[4,5]decane-7-etha...	212	C12H20O3	020489-42-3	28



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG041515\
Data File : BG016439.D
Acq On : 16 Apr 2015 2:33
Operator : TP/IZ
Sample : G1829-12
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
TE-PMW-1A-11

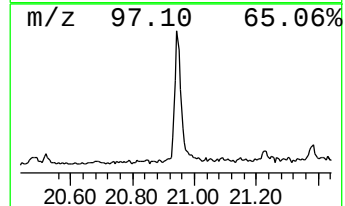
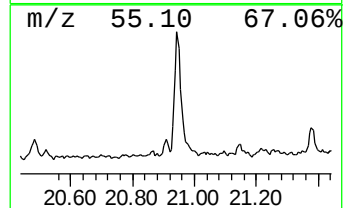
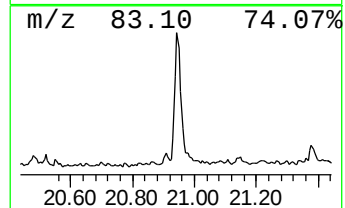
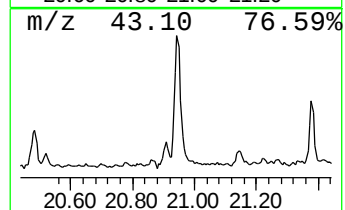
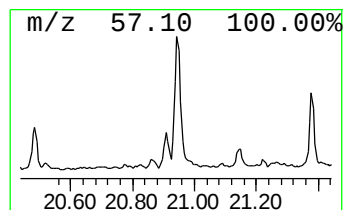
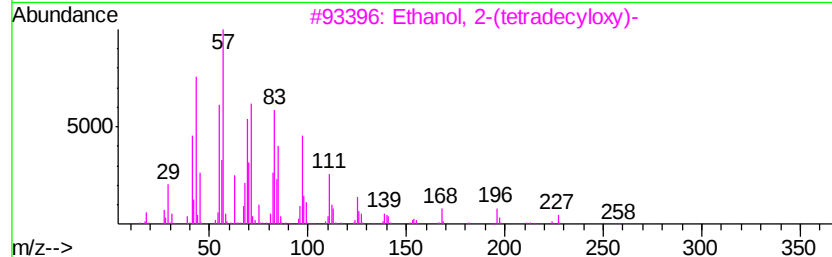
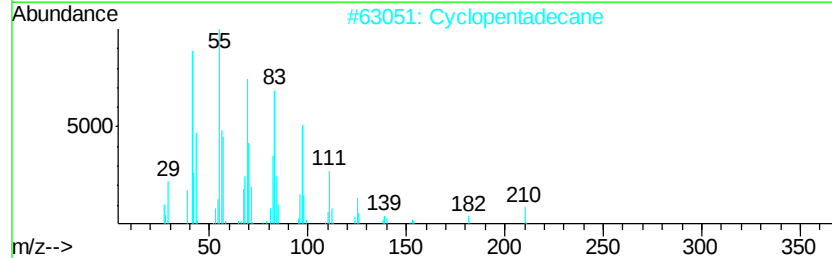
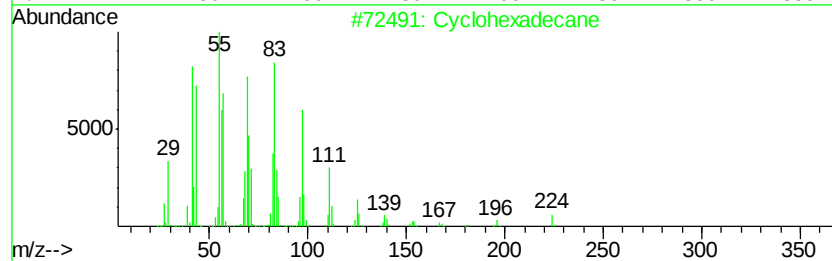
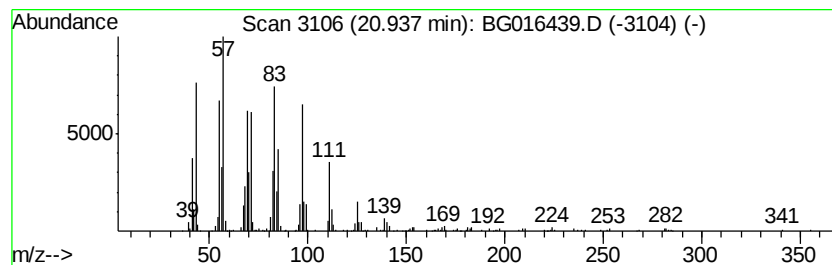
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Cyclohexadecane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.94	8.15 ng	502726	Chrysene-d12	21.23

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclohexadecane	224	C16H32	000295-65-8	93
2		Cyclopentadecane	210	C15H30	000295-48-7	92
3		Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	91
4		1-Docosene	308	C22H44	001599-67-3	90
5		Heptafluorobutanoic acid, heptad...	452	C21H35F7O2	1000282-97-3	87



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG041515\
Data File : BG016439.D
Acq On : 16 Apr 2015 2:33
Operator : TP/IZ
Sample : G1829-12
Misc :
ALS Vial : 4 Sample Multiplier: 1

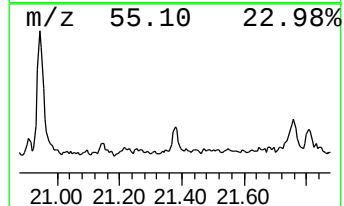
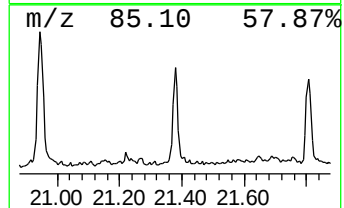
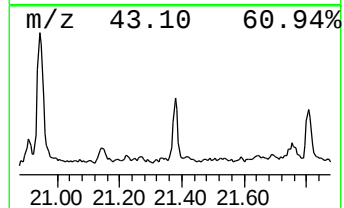
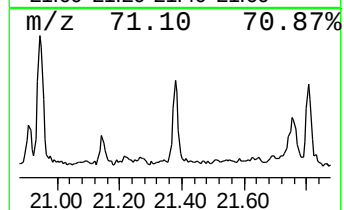
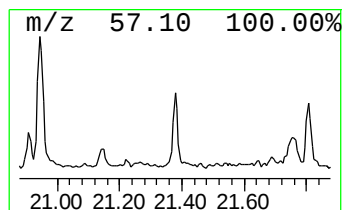
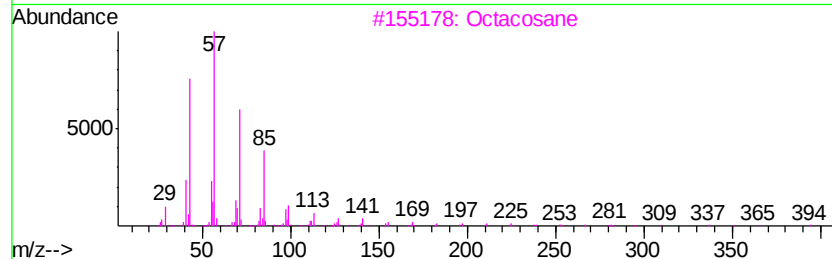
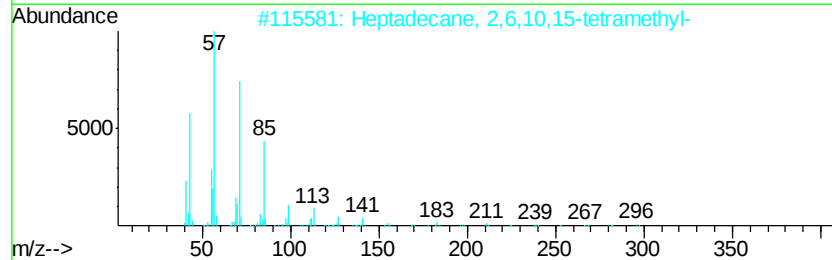
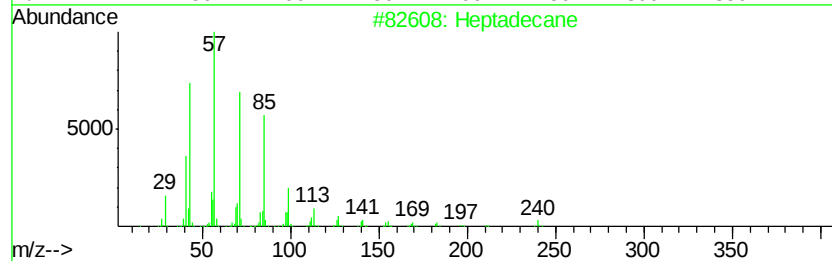
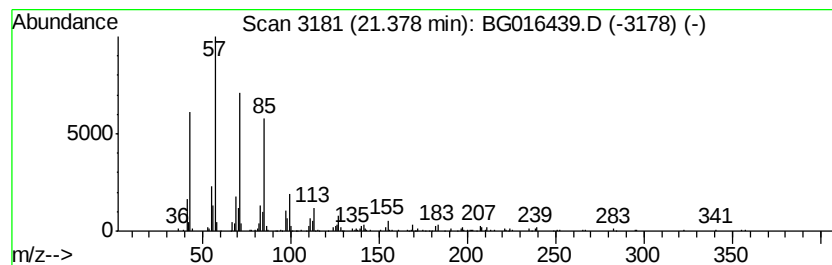
Instrument :
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG040615.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Heptadecane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.38	2.03 ng	125456	Chrysene-d12	21.23
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Heptadecane	240 C17H36	000629-78-7	91
2	Heptadecane, 2,6,10,15-tetramethyl-	296 C21H44	054833-48-6	91
3	Octacosane	394 C28H58	000630-02-4	91
4	Heneicosane	296 C21H44	000629-94-7	91
5	Heptacosane	380 C27H56	000593-49-7	91



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG041515\
Data File : BG016439.D
Acq On : 16 Apr 2015 2:33
Operator : TP/IZ
Sample : G1829-12
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
TE-PMW-1A-11

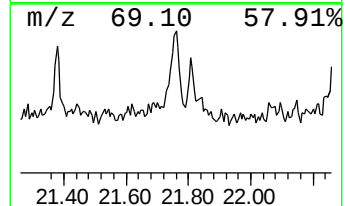
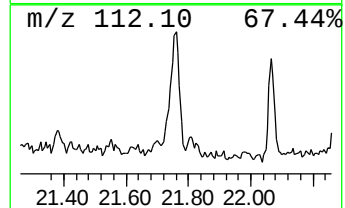
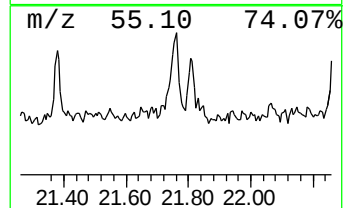
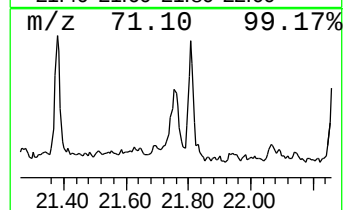
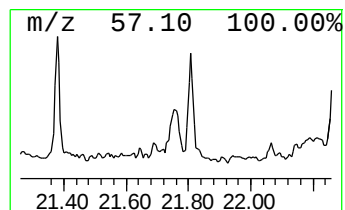
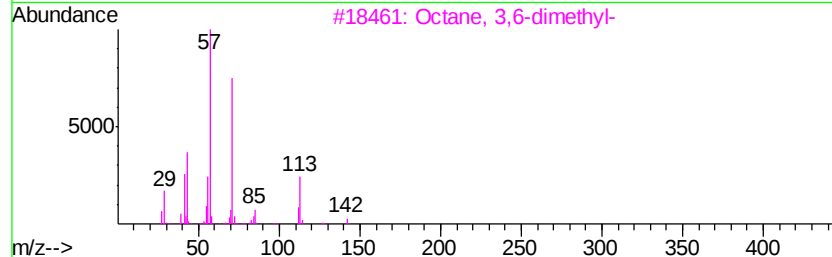
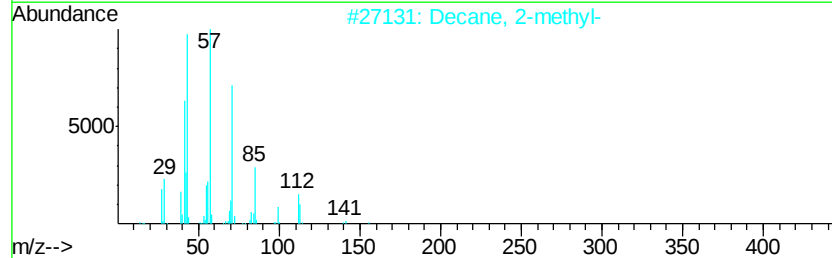
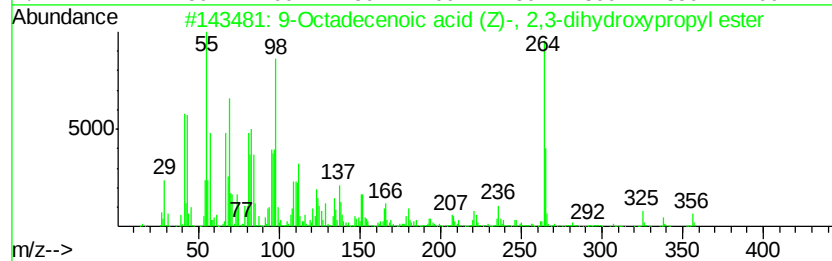
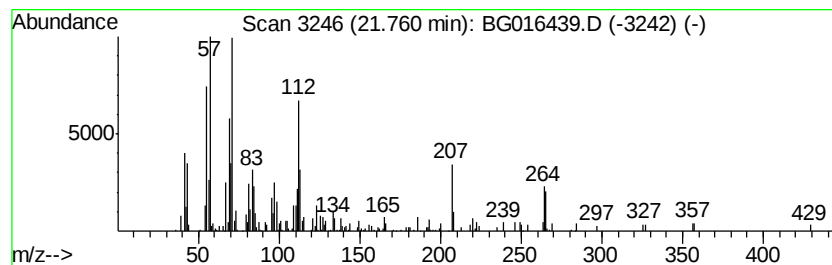
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 unknown21.76 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.76	2.22 ng	137286	Chrysene-d12	21.23

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9-Octadecenoic acid (Z)-, 2,3-di...	356	C21H40O4	000111-03-5	41
2		Decane, 2-methyl-	156	C11H24	006975-98-0	35
3		Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	27
4		2-Bromo-6-methylheptane	192	C8H17Br	004730-24-9	22
5		3-Cyclopentylpropionic acid, 2-t...	226	C13H22O3	1000293-46-9	22



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG041515\
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Acq On : 16 Apr 2015 2:33
Operator : TP/IZ
Sample : G1829-12
Misc :
ALS Vial : 4 Sample Multiplier: 1

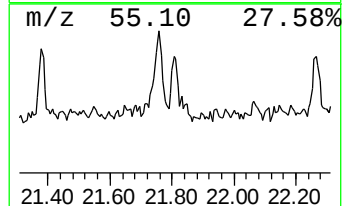
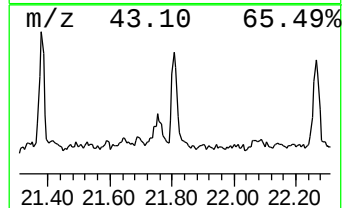
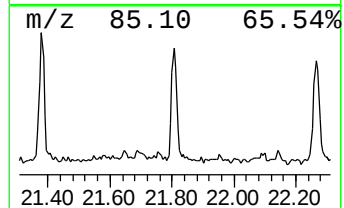
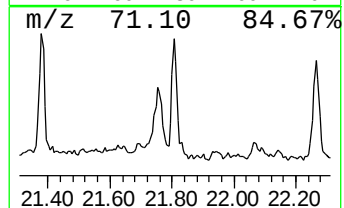
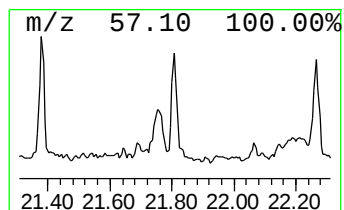
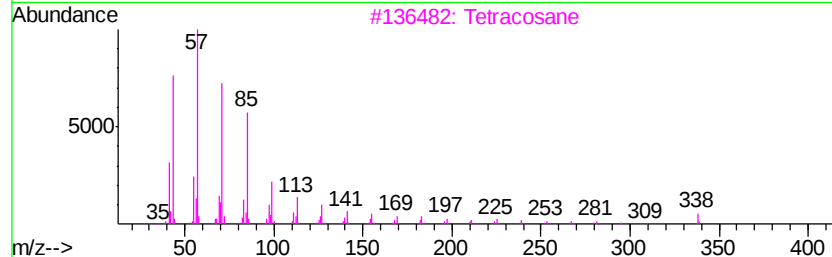
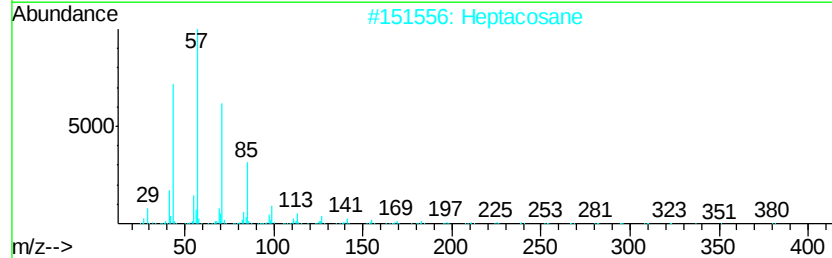
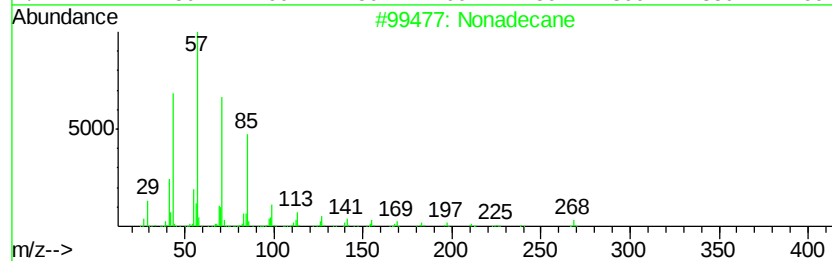
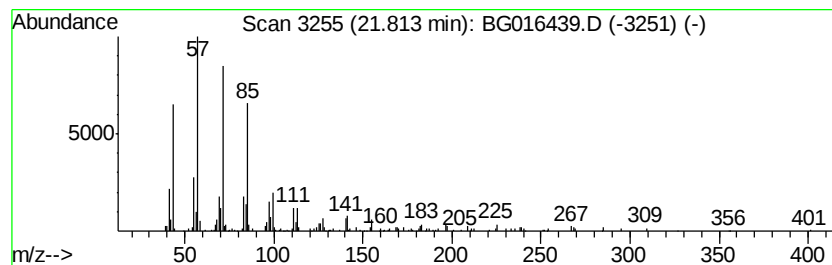
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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG040615.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Nonadecane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.81	2.20 ng	135939	Chrysene-d12	21.23
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Nonadecane	268 C19H40	000629-92-5	95
2	Heptacosane	380 C27H56	000593-49-7	94
3	Tetracosane	338 C24H50	000646-31-1	91
4	Heneicosane	296 C21H44	000629-94-7	91
5	Tetratetracontane	619 C44H90	007098-22-8	91



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG041515\
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Acq On : 16 Apr 2015 2:33
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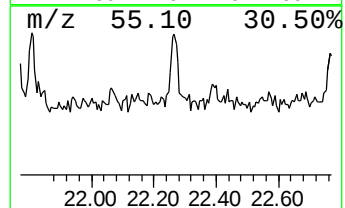
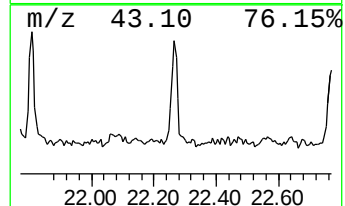
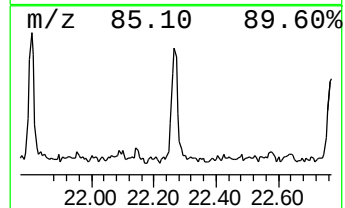
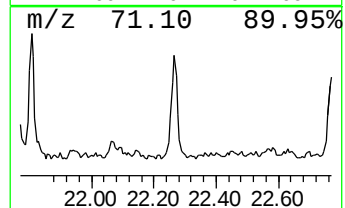
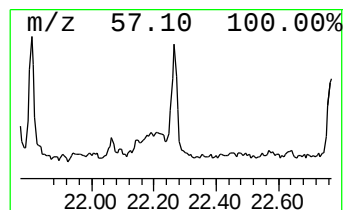
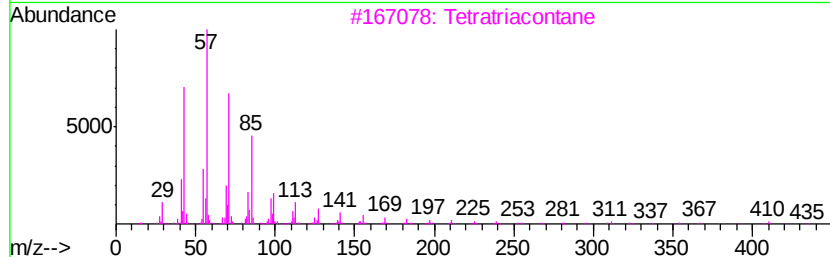
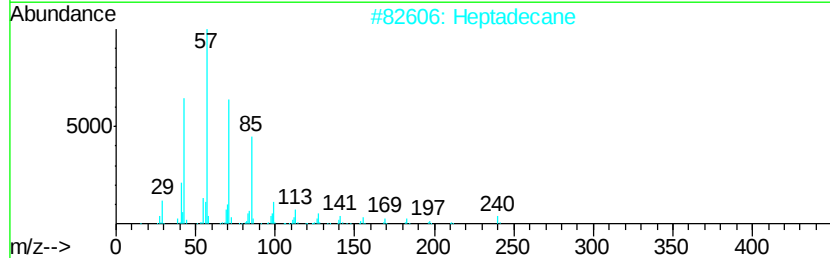
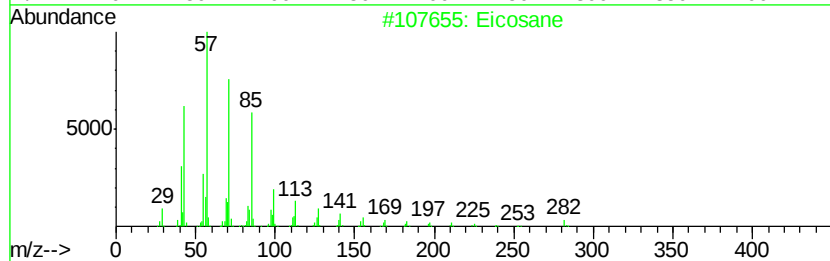
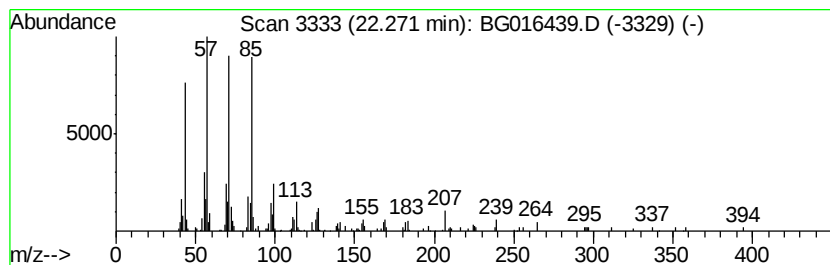
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG040615.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Eicosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.27	2.63 ng	162486	Chrysene-d12	21.23
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Eicosane		282 C20H42	000112-95-8 97
2	Heptadecane		240 C17H36	000629-78-7 94
3	Tetratriacontane		479 C34H70	014167-59-0 91
4	Nonadecane, 9-methyl-		282 C20H42	013287-24-6 91
5	Heneicosane		296 C21H44	000629-94-7 91



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG041515\
Data File : BG016439.D
Acq On : 16 Apr 2015 2:33
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Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
TE-PMW-1A-11

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG040615.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
Butane, 2-methoxy...	2.83	20.0	ng	450906	1	7.66	450803	20.0
2-Pentanone, 4-hy...	4.78	4.1	ng	93321	1	7.66	450803	20.0
unknown7.20	7.20	88.3	ng	1989800	1	7.66	450803	20.0
n-Hexadecanoic acid	17.95	4.3	ng	278565	4	17.05	1288890	20.0
Octadecanoic acid	19.23	2.3	ng	144595	5	21.23	1234040	20.0
1,2-Cyclohexanedi...	20.91	2.2	ng	133971	5	21.23	1234040	20.0
Cyclohexadecane	20.94	8.2	ng	502726	5	21.23	1234040	20.0
Heptadecane	21.38	2.0	ng	125456	5	21.23	1234040	20.0
unknown21.76	21.76	2.2	ng	137286	5	21.23	1234040	20.0
Nonadecane	21.81	2.2	ng	135939	5	21.23	1234040	20.0
Eicosane	22.27	2.6	ng	162486	5	21.23	1234040	20.0