

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110921\
 Data File : BP007903.D
 Acq On : 09 Nov 2021 22:37
 Operator : CG/JU
 Sample : M4485-07
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 AG-8

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:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP110221.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP007903.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.446	360	367	377	rBV	1635922	2342907	43.46%	8.938%
2	7.034	629	637	650	rBV	1475309	2344269	43.48%	8.943%
3	7.858	770	777	785	rBV	443660	697095	12.93%	2.659%
4	9.016	966	974	985	rBV	798086	1349936	25.04%	5.150%
5	10.446	1207	1217	1228	rBV	86509	166919	3.10%	0.637%
6	10.657	1246	1253	1264	rBV	486055	831208	15.42%	3.171%
7	13.116	1664	1671	1691	rBV	2149561	3241339	60.12%	12.365%
8	14.492	1899	1905	1919	rBV	723750	1091664	20.25%	4.165%
9	15.987	2152	2159	2180	rBV	1793282	2659900	49.34%	10.147%
10	17.251	2367	2374	2387	rBV	883947	1347637	25.00%	5.141%
11	18.145	2521	2526	2532	rBV2	51549	80289	1.49%	0.306%
12	19.810	2804	2809	2819	rBV	4305900	5391343	100.00%	20.567%
13	21.051	3017	3020	3027	rBV	302189	427514	7.93%	1.631%
14	21.339	3064	3069	3077	rBV	1431647	1837389	34.08%	7.009%
15	21.980	3174	3178	3189	rBV	252505	362991	6.73%	1.385%
16	23.751	3473	3479	3489	rVB	996803	2040817	37.85%	7.785%

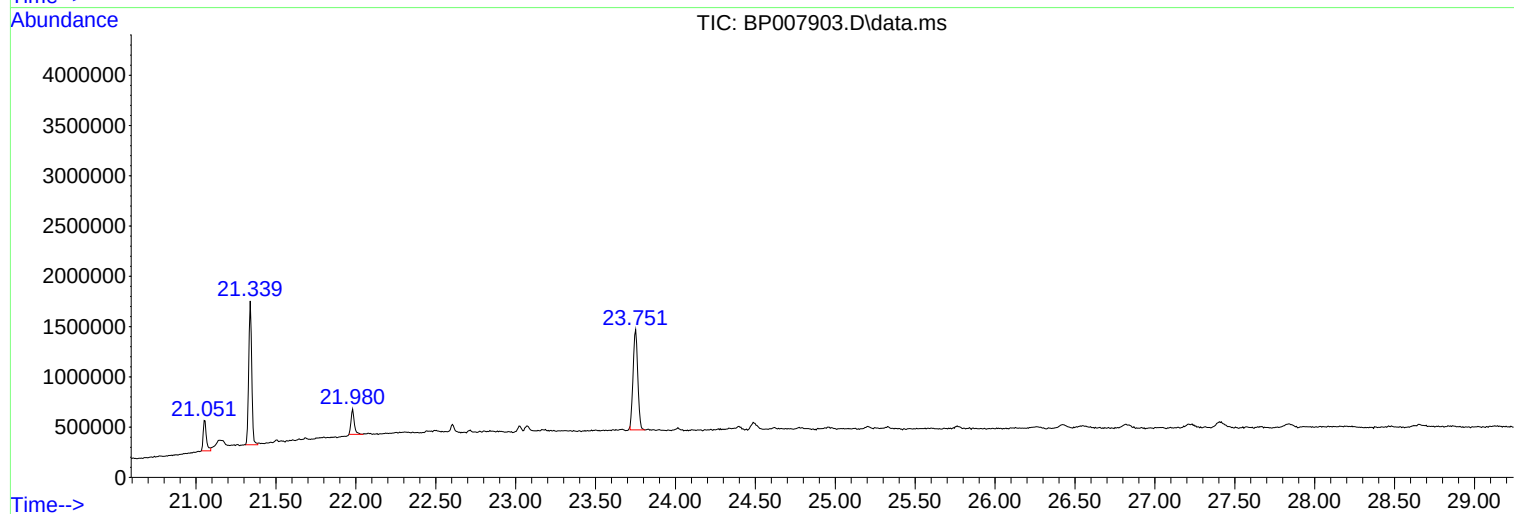
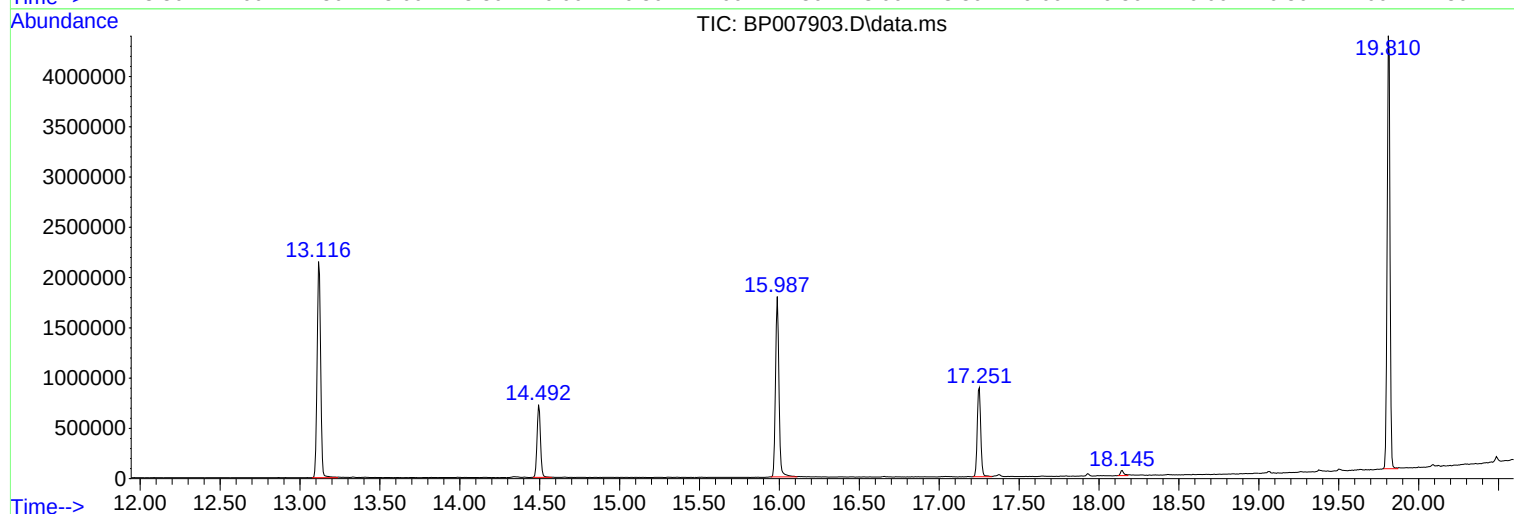
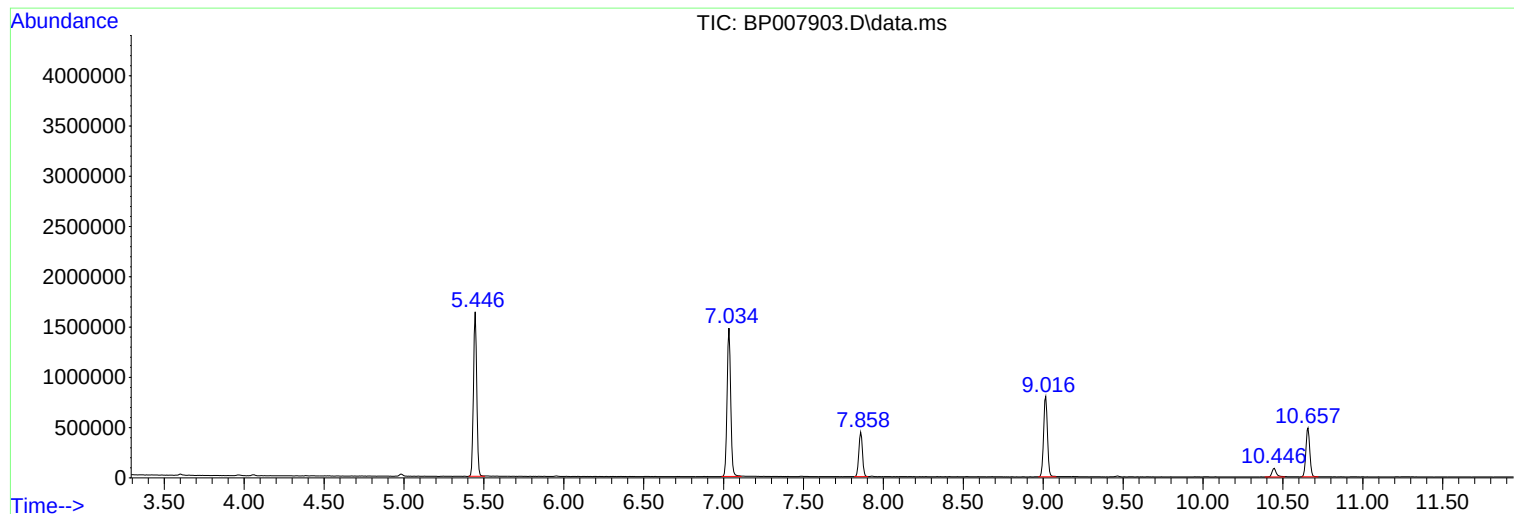
Sum of corrected areas: 26213217

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP110221.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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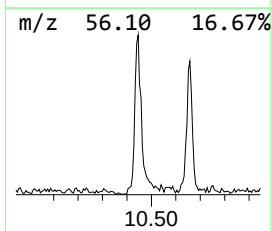
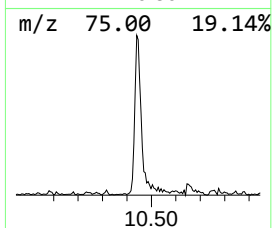
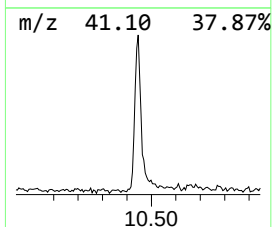
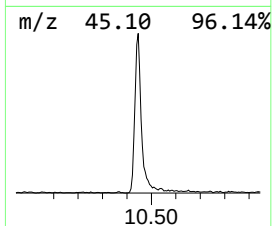
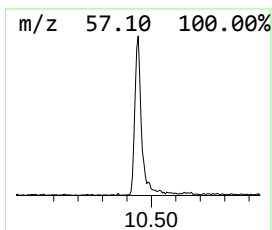
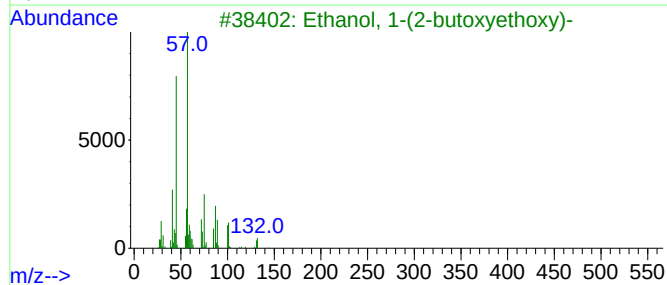
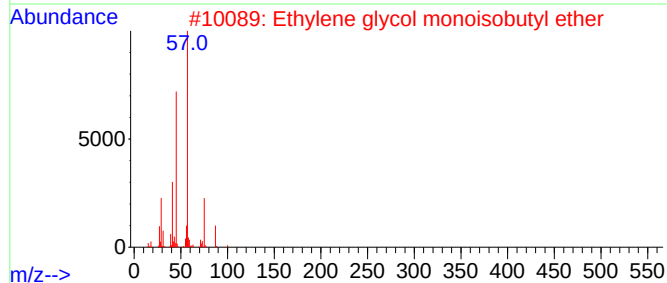
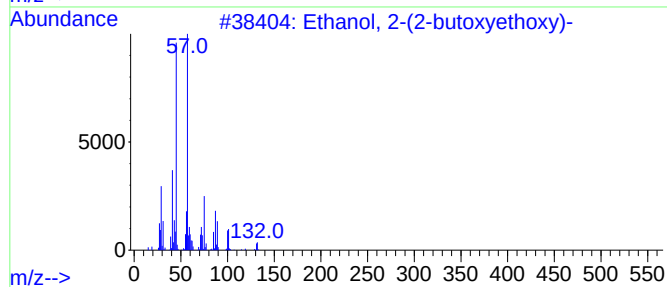
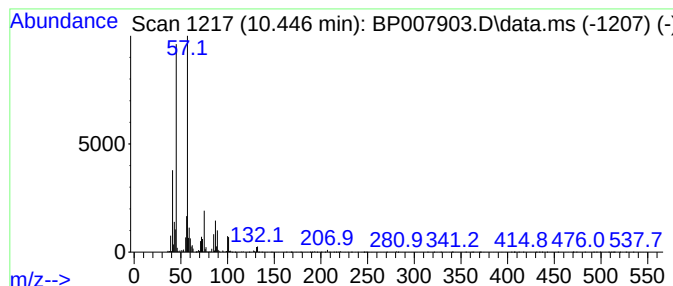
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP110221.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.446	4.02 ng	166919	Naphthalene-d8	10.657

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
2			Ethylene glycol monoisobutyl ether	118	C6H14O2	004439-24-1	64
3			Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	58
4			3,6,9,12-Tetraoxahexadecan-1-ol	250	C12H26O5	001559-34-8	53
5			Hexaethylene glycol monododecyl ...	450	C24H50O7	003055-96-7	50



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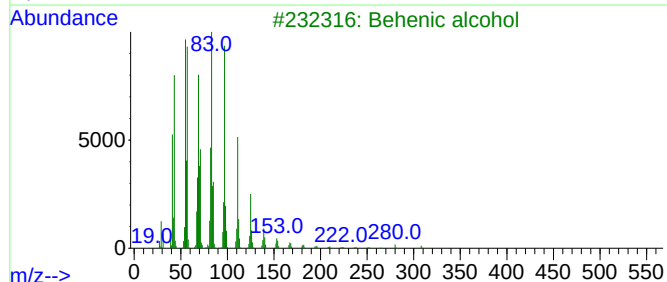
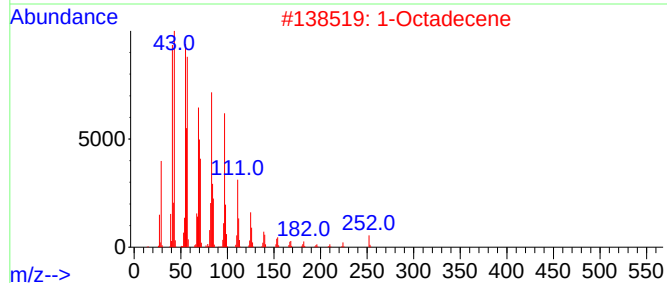
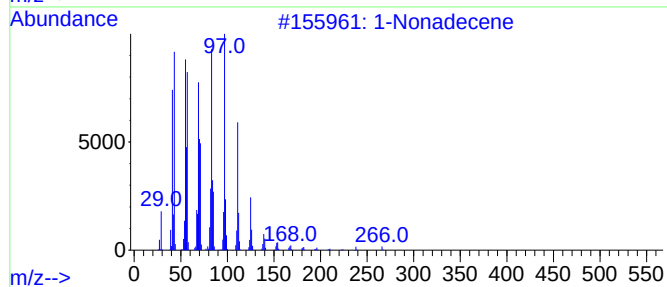
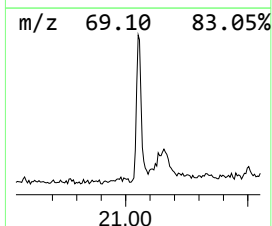
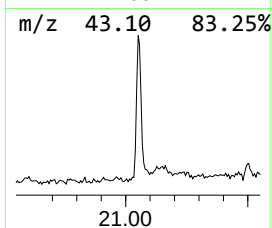
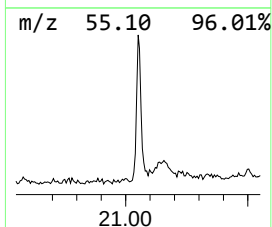
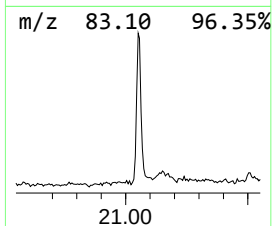
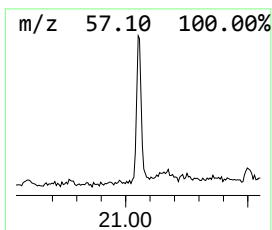
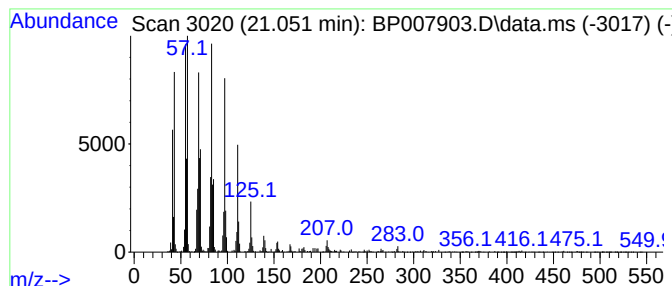
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Peak Number 2 1-Nonadecene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.051	4.65 ng	427514	Chrysene-d12	21.339

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Nonadecene			266	C19H38	018435-45-5	98
2	1-Octadecene			252	C18H36	000112-88-9	96
3	Behenic alcohol			326	C22H46O	000661-19-8	95
4	n-Tetracosanol-1			354	C24H50O	000506-51-4	95
5	1-Tetracosene			336	C24H48	010192-32-2	95



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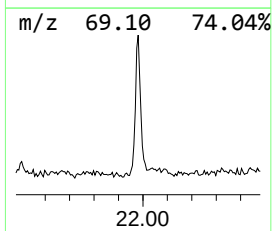
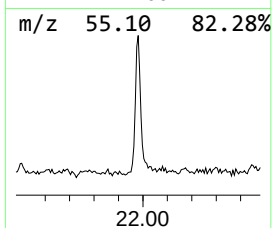
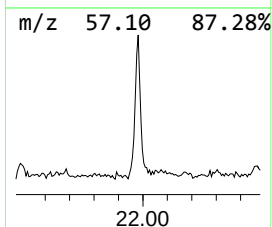
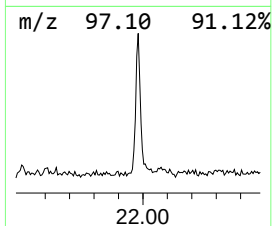
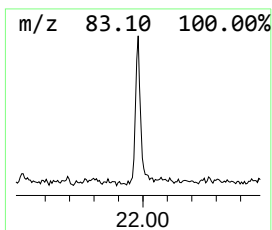
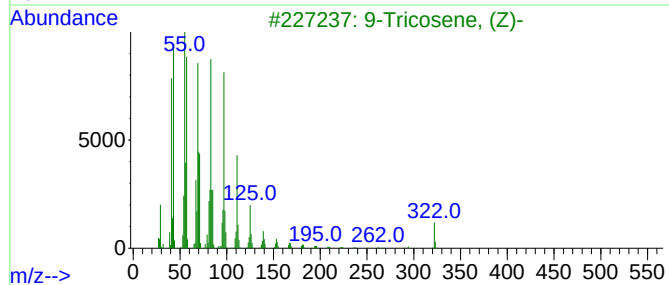
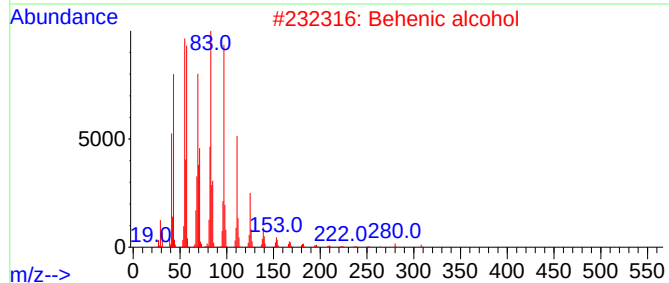
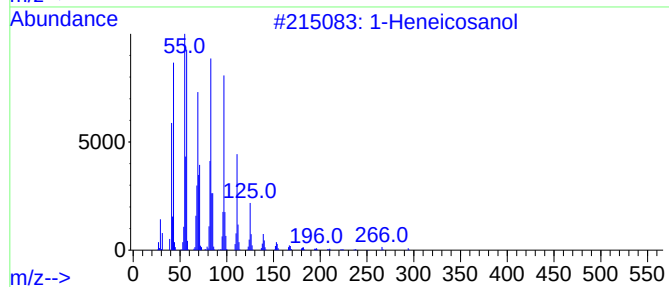
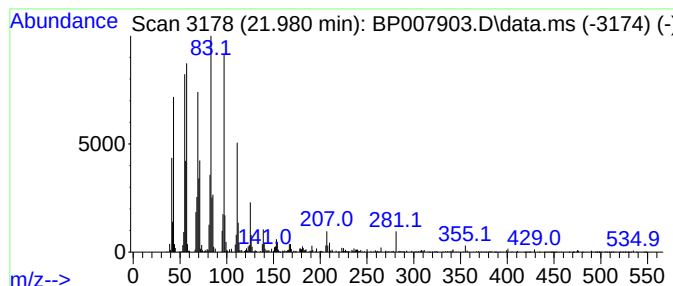
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Peak Number 3 1-Heneicosanol Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.980	3.95 ng	362991	Chrysene-d12	21.339

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Heneicosanol	312	C21H44O	015594-90-8	95
2			Behenic alcohol	326	C22H46O	000661-19-8	95
3			9-Tricosene, (Z)-	322	C23H46	027519-02-4	94
4			1-Docosanol, methyl ether	340	C23H48O	1000333-91-9	94
5			n-Nonadecanol-1	284	C19H40O	001454-84-8	94



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TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Ethanol, 2-(2-b...	10.446	4.0	ng	166919	2	10.657	831208	20.0
1-Nonadecene	21.051	4.7	ng	427514	5	21.339	1837390	20.0
1-Heneicosanol	21.980	4.0	ng	362991	5	21.339	1837390	20.0