

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121518\
 Data File : BM018212.D
 Acq On : 15 Dec 2018 20:08
 Operator : JU/SJ
 Sample : J6347-06
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 P001-SW001-01

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_M\METHODS\8270-BM121518.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	5.128	376	381	389	rBV	92254	147233	7.47%	0.689%
2	6.704	641	649	650	rBV2	63749	100777	5.12%	0.472%
3	6.722	650	652	661	rVB	84084	136582	6.93%	0.639%
4	7.040	699	706	716	rBV	170159	287658	14.60%	1.347%
5	7.128	716	721	735	rVV	66578	153859	7.81%	0.720%
6	7.498	777	784	792	rBV	455938	723322	36.72%	3.387%
7	7.810	830	837	847	rBV	222725	359308	18.24%	1.682%
8	8.663	973	982	991	rBV	131716	267805	13.59%	1.254%
9	9.034	1040	1045	1051	rVB2	21121	30671	1.56%	0.144%
10	10.269	1248	1255	1275	rBV	540385	1033417	52.46%	4.838%
11	12.769	1672	1680	1696	rBV	455666	779420	39.56%	3.649%
12	13.645	1823	1829	1837	rBV2	33307	68236	3.46%	0.319%
13	14.157	1910	1916	1930	rBV2	933334	1434850	72.84%	6.718%
14	15.669	2164	2173	2181	rBV	303789	504991	25.63%	2.364%
15	16.698	2342	2348	2356	rBV7	8729	21908	1.11%	0.103%
16	16.921	2380	2386	2393	rBV	1113456	1677348	85.15%	7.853%
17	17.221	2431	2437	2443	rBV	70629	100372	5.10%	0.470%
18	17.704	2512	2519	2527	rBV5	9613	24065	1.22%	0.113%
19	17.833	2536	2541	2551	rBV	713206	1055292	53.57%	4.941%
20	18.980	2731	2736	2738	rBV4	29261	43967	2.23%	0.206%
21	19.127	2756	2761	2774	rBV2	389729	586580	29.78%	2.746%
22	19.574	2831	2837	2845	rBV	1114183	1443014	73.25%	6.756%
23	19.868	2884	2887	2889	rBV3	24404	25610	1.30%	0.120%
24	19.892	2889	2891	2894	rVB3	26920	23003	1.17%	0.108%
25	20.392	2973	2976	2980	rVB	67759	69327	3.52%	0.325%
26	20.498	2992	2994	2999	rBV6	15523	22521	1.14%	0.105%
27	20.862	3051	3056	3063	rBV	786243	1005912	51.06%	4.710%
28	21.062	3086	3090	3094	rBV	210153	227406	11.54%	1.065%
29	21.139	3098	3103	3112	rBV	1446975	1969984	100.00%	9.223%
30	21.298	3126	3130	3136	rBV	411872	469593	23.84%	2.199%
31	21.721	3198	3202	3208	rBV	595852	723004	36.70%	3.385%
32	22.168	3273	3278	3286	rVB	791553	962941	48.88%	4.508%
33	22.280	3295	3297	3301	rBV2	57729	63009	3.20%	0.295%
34	22.650	3355	3360	3366	rBV	814243	1038749	52.73%	4.863%

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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	23.186	3446	3451	3458	rVB	649395	1033151	52.44%	4.837%
36	23.303	3464	3471	3481	rVB	800300	1614821	81.97%	7.560%
37	23.791	3549	3554	3560	rVB	403462	714730	36.28%	3.346%
38	24.497	3668	3674	3680	rVB2	222592	414322	21.03%	1.940%

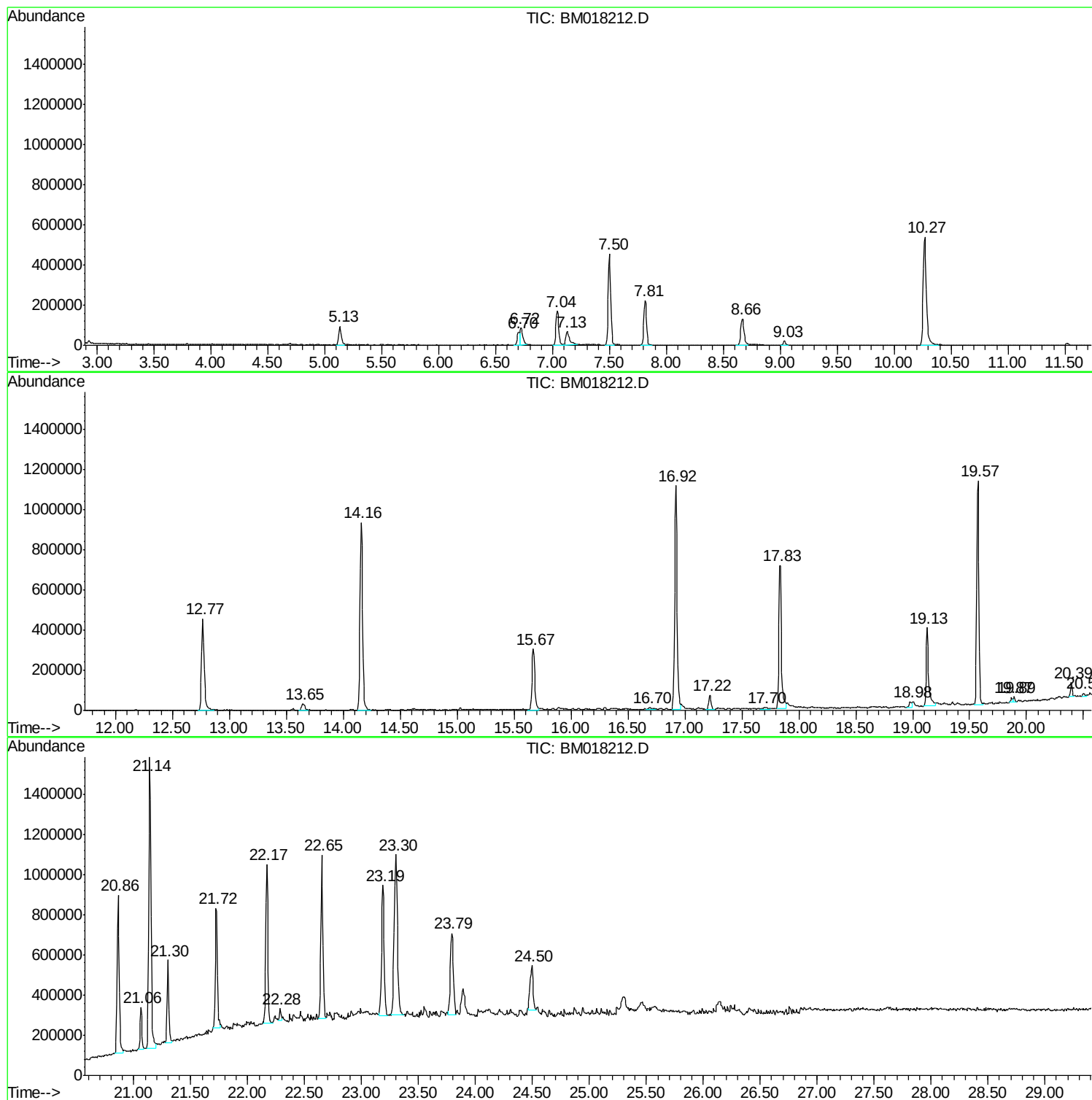
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\8270-BM121518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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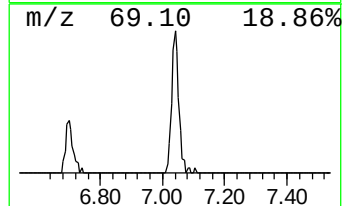
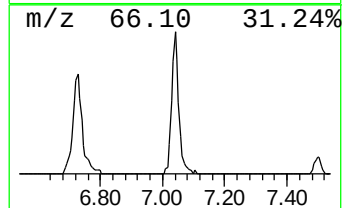
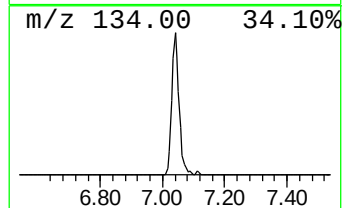
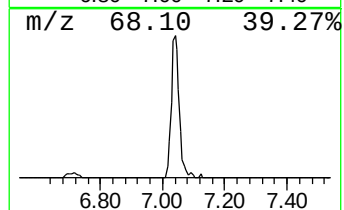
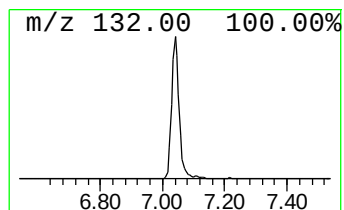
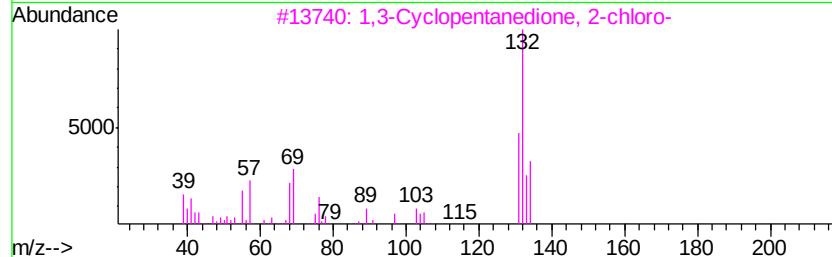
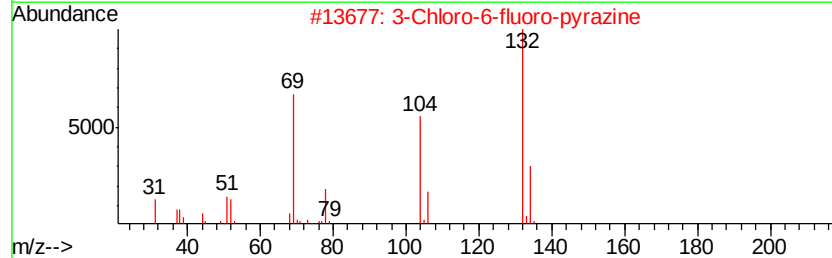
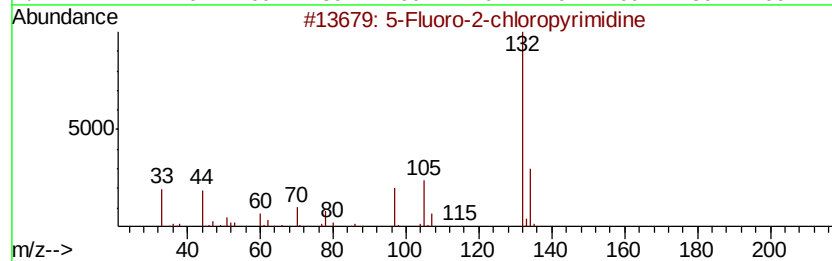
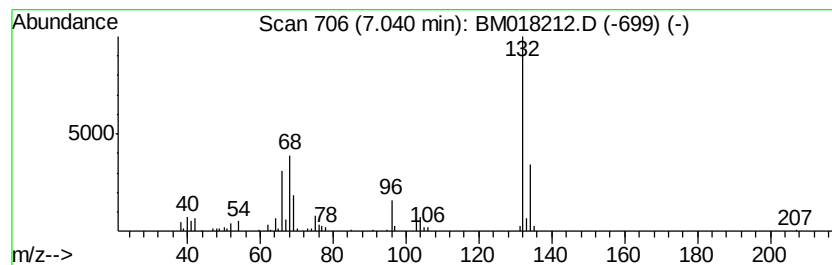
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\8270-BM121518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown7.04 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.04	7.95 ng	287658	1,4-Dichlorobenzene-d4	7.50

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	27
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
3		1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	25
4		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
5		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12



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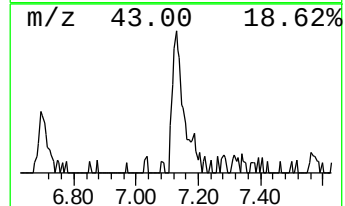
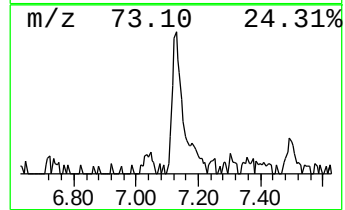
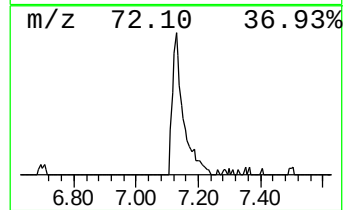
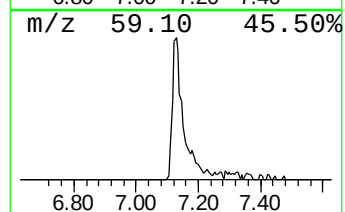
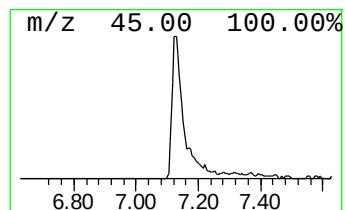
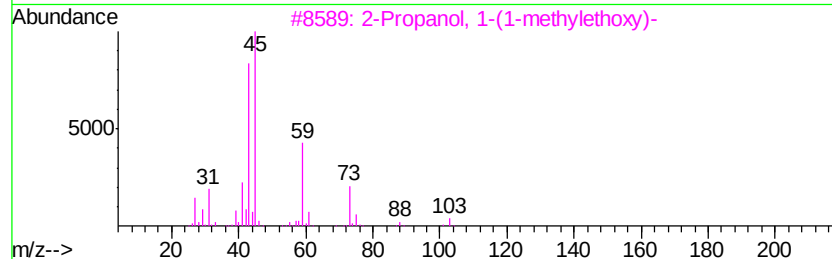
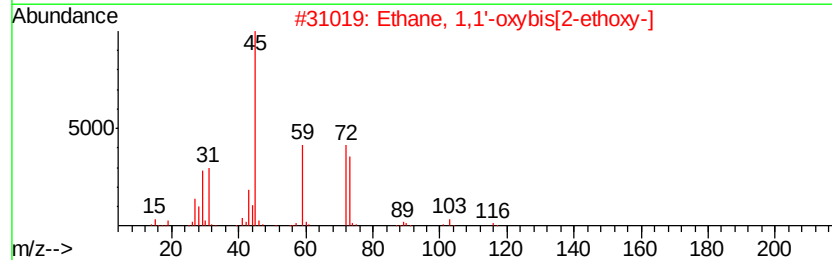
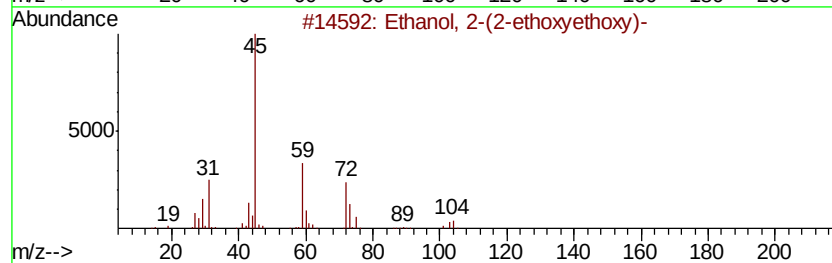
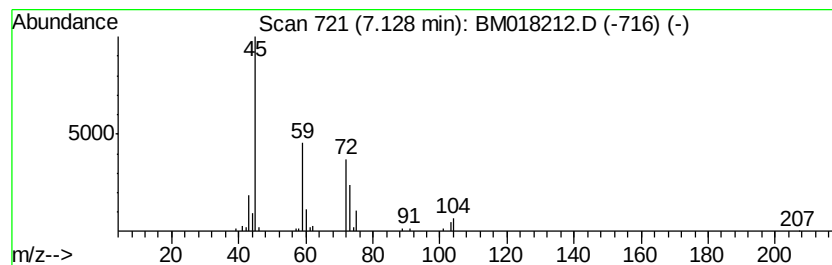
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\8270-BM121518.M
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 2 Ethanol, 2-(2-ethoxyethoxy)- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.13	4.25 ng	153859	1,4-Dichlorobenzene-d4	7.50

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Ethanol, 2-(2-ethoxyethoxy)-	134	C6H14O3	000111-90-0	90
2		Ethane, 1,1'-oxybis[2-ethoxy-]	162	C8H18O3	000112-36-7	72
3		2-Propanol, 1-(1-methylethoxy)-	118	C6H14O2	003944-36-3	50
4		Ethanol, 2-[2-(2-ethoxyethoxy)et...]	178	C8H18O4	000112-50-5	45
5		Ethanol, 2-(2-methoxyethoxy)-	120	C5H12O3	000111-77-3	38



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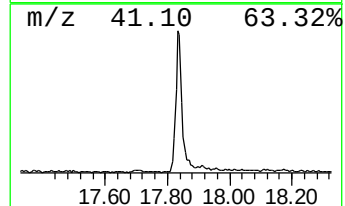
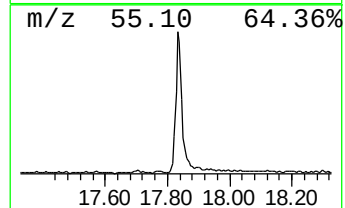
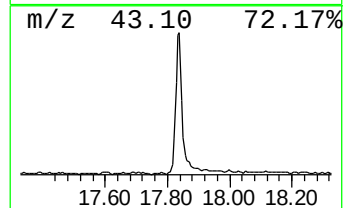
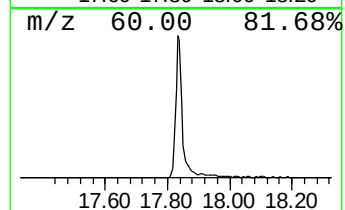
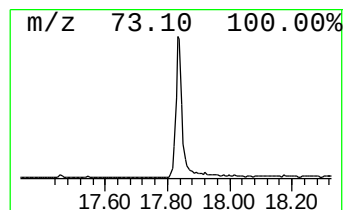
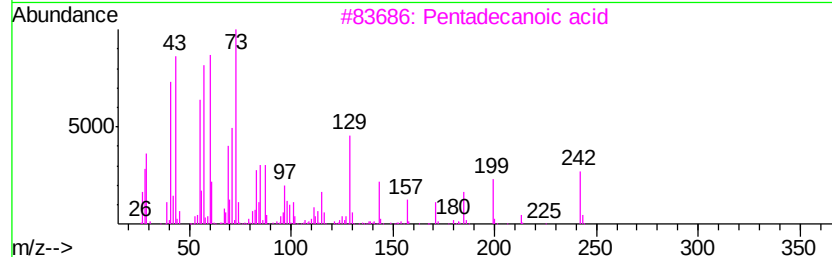
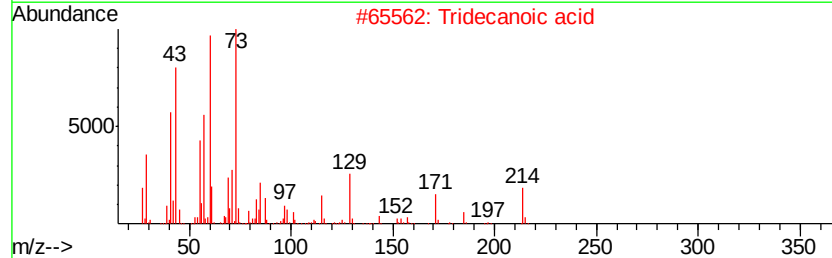
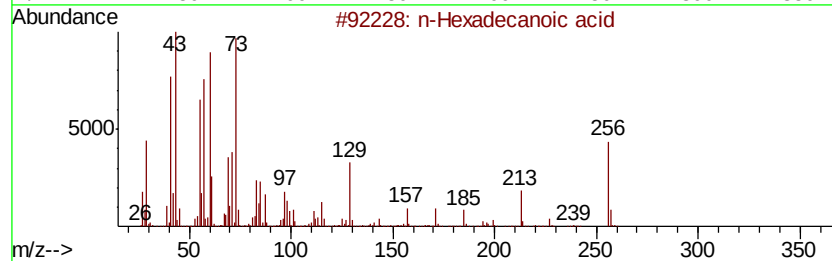
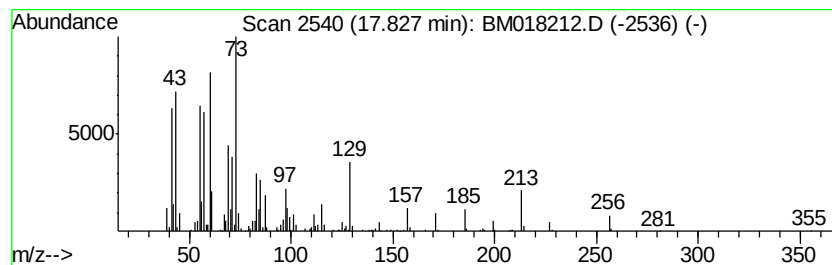
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD		R.T.	
17.83	12.58 ng	1055290	Phenanthrene-d10		16.92	
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	89
3		Pentadecanoic acid	242	C15H30O2	001002-84-2	74
4		Ether, dodecyl isopropyl	228	C15H32O	029379-42-8	38
5		Dodecane	170	C12H26	000112-40-3	11



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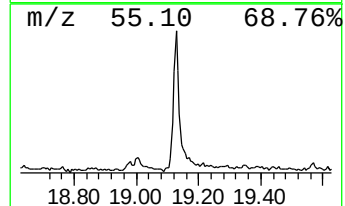
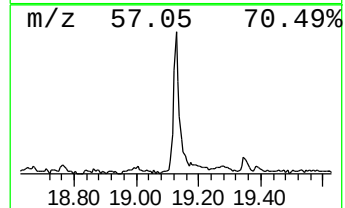
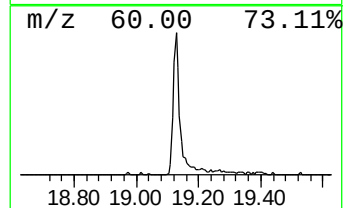
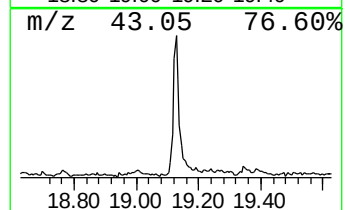
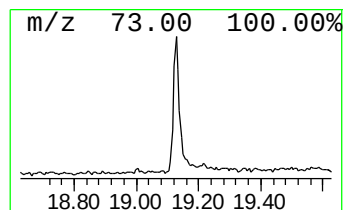
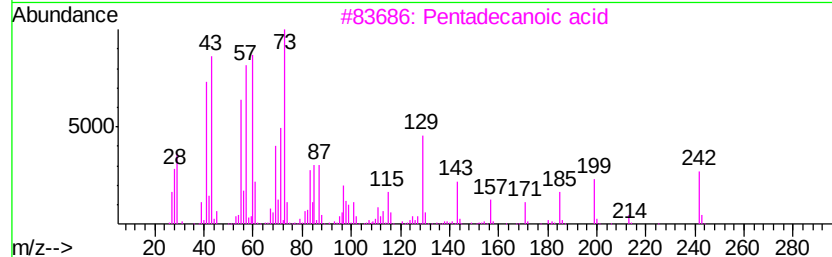
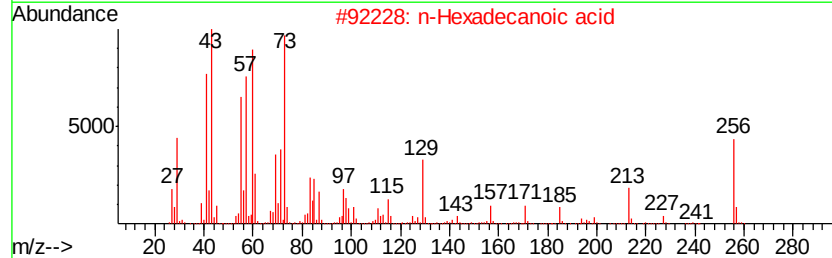
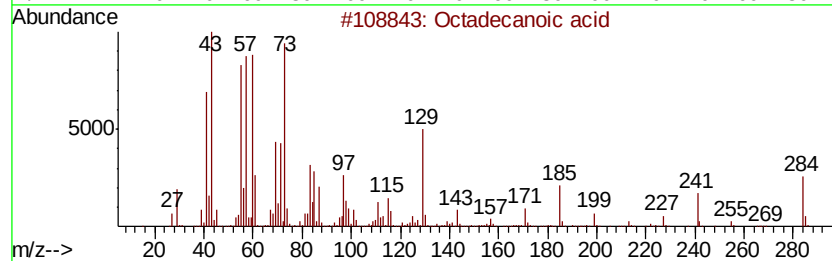
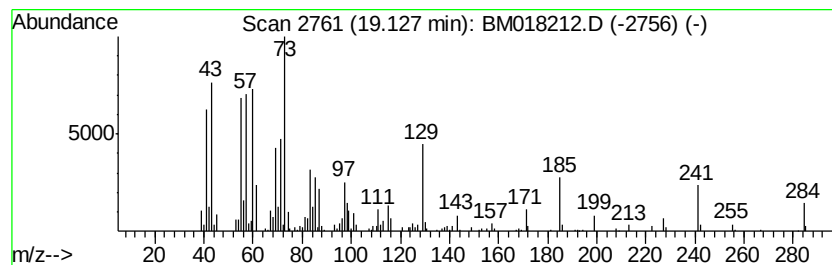
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Octadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.13	5.96 ng	586580	Chrysene-d12	21.14

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	93
3		Pentadecanoic acid	242	C15H30O2	001002-84-2	76
4		n-Decanoic acid	172	C10H20O2	000334-48-5	64
5		Tetradecanoic acid	228	C14H28O2	000544-63-8	58



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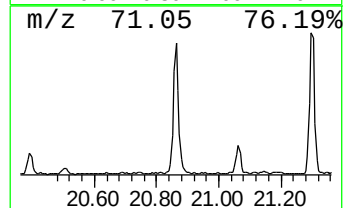
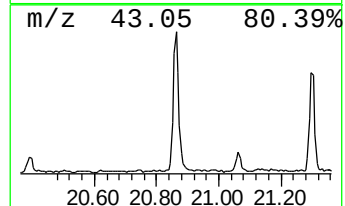
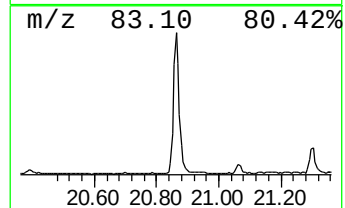
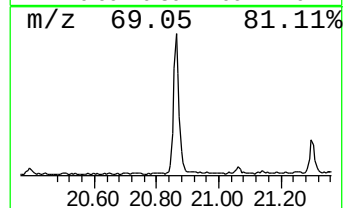
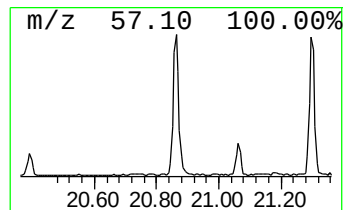
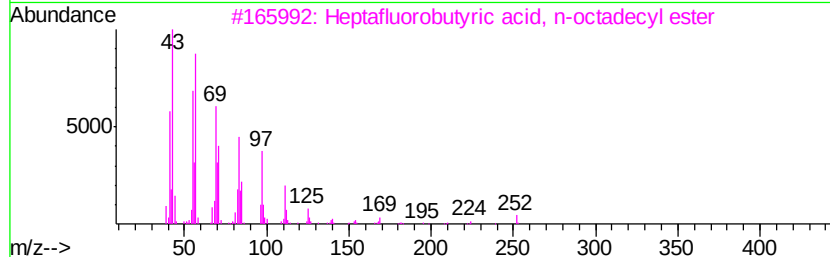
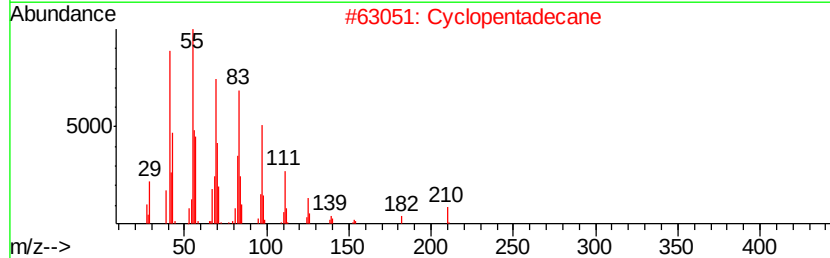
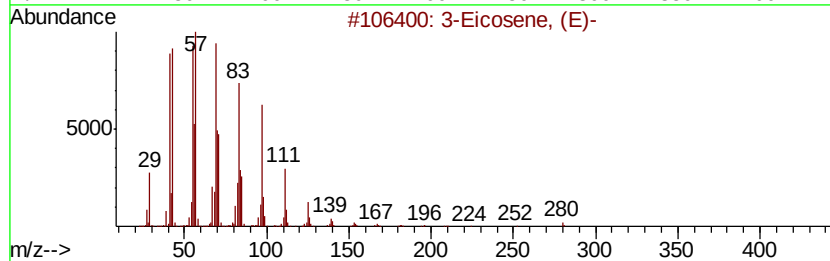
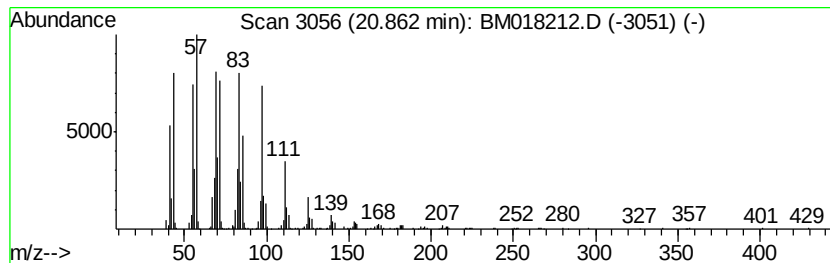
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 3-Eicosene, (E)- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.86	10.21 ng	1005910	Chrysene-d12	21.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Eicosene, (E)-	280	C20H40	074685-33-9	96
2		Cyclopentadecane	210	C15H30	000295-48-7	87
3		Heptafluorobutyric acid, n-octad...	466	C22H37F7O2	000400-57-7	87
4		Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	81
5		Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	74



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121518\
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Acq On : 15 Dec 2018 20:08
Operator : JU/SJ
Sample : J6347-06
Misc :
ALS Vial : 13 Sample Multiplier: 1

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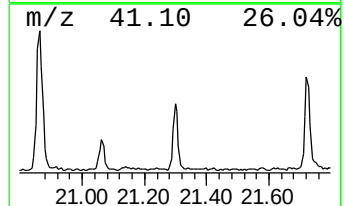
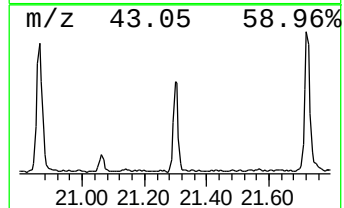
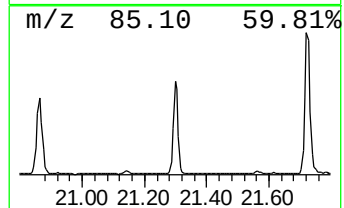
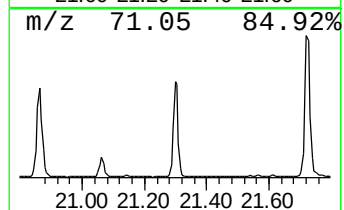
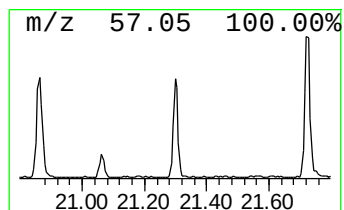
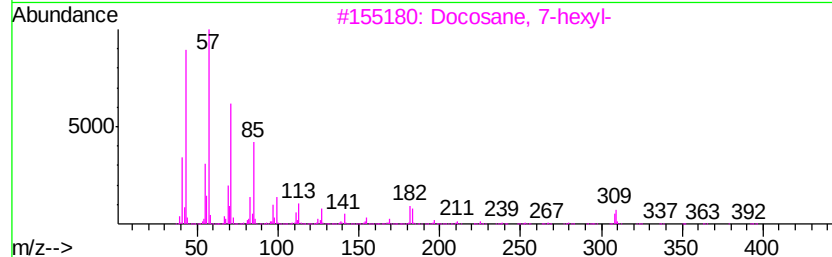
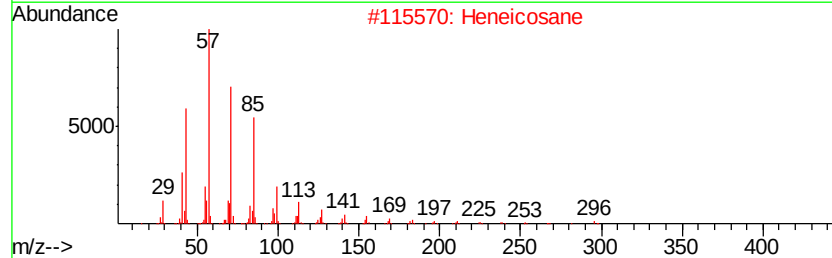
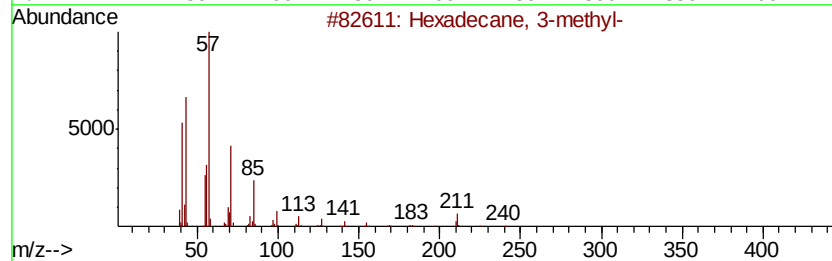
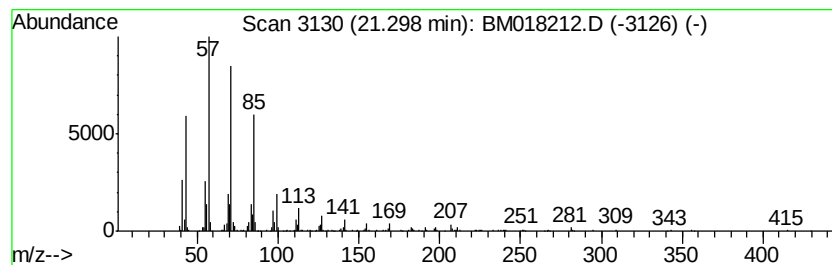
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Hexadecane, 3-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.30	4.77 ng	469593	Chrysene-d12	21.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecane, 3-methyl-	240	C17H36	006418-43-5	91
2		Heneicosane	296	C21H44	000629-94-7	91
3		Docosane, 7-hexyl-	394	C28H58	055373-86-9	91
4		Docosane, 11-butyl-	366	C26H54	013475-76-8	91
5		Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121518\
Data File : BM018212.D
Acq On : 15 Dec 2018 20:08
Operator : JU/SJ
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Misc :
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ClientSampleId :
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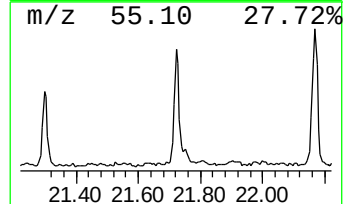
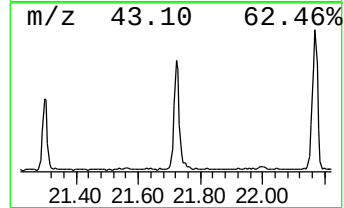
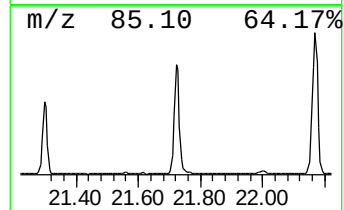
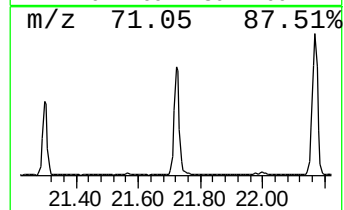
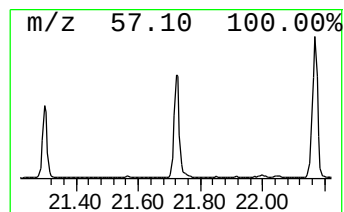
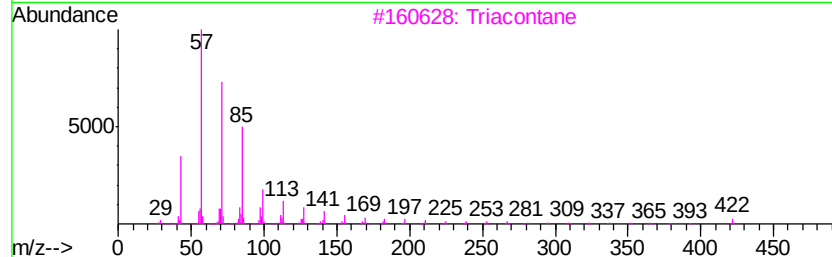
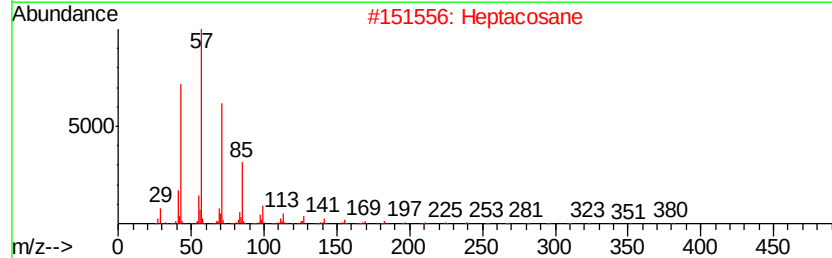
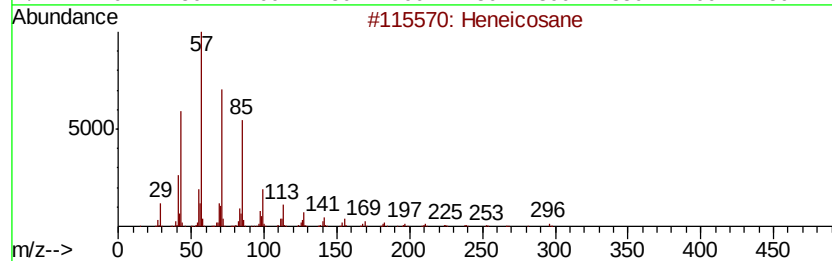
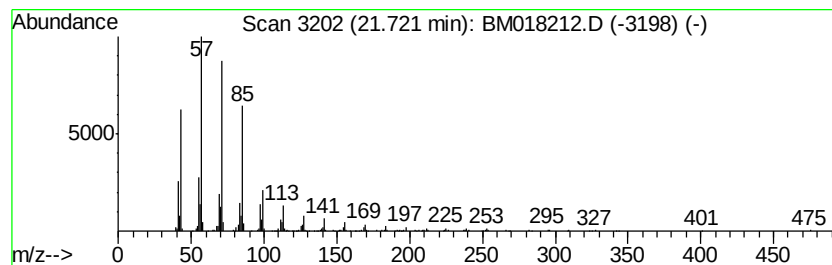
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Heneicosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.72	7.34 ng	723004	Chrysene-d12	21.14

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane		296	C21H44	000629-94-7	91
2	Heptacosane		380	C27H56	000593-49-7	91
3	Triacontane		422	C30H62	000638-68-6	91
4	Tridecanol, 2-ethyl-2-methyl-		242	C16H34O	1000115-66-1	90
5	Octacosane		394	C28H58	000630-02-4	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121518\
Data File : BM018212.D
Acq On : 15 Dec 2018 20:08
Operator : JU/SJ
Sample : J6347-06
Misc :
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Instrument :
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ClientSampleId :
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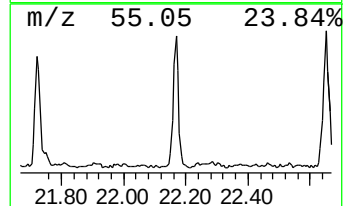
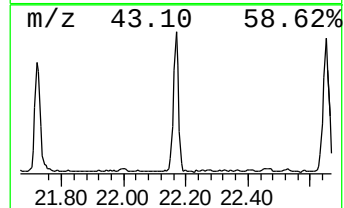
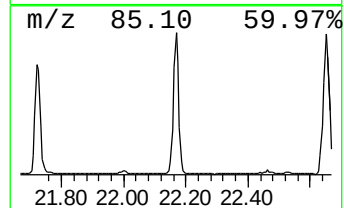
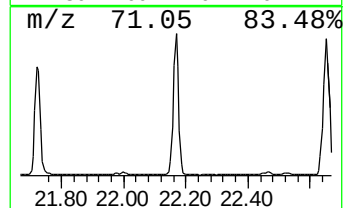
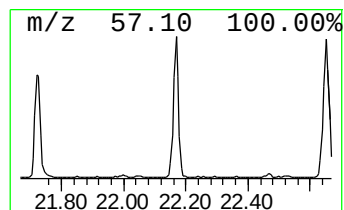
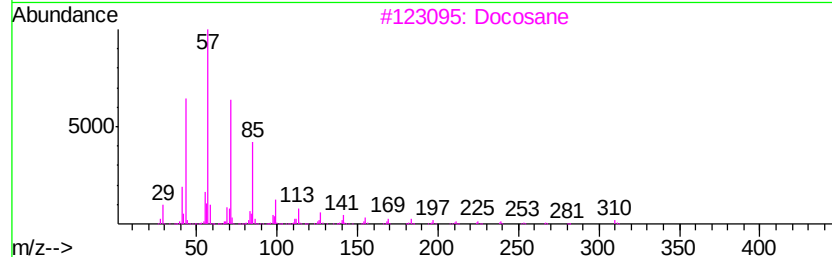
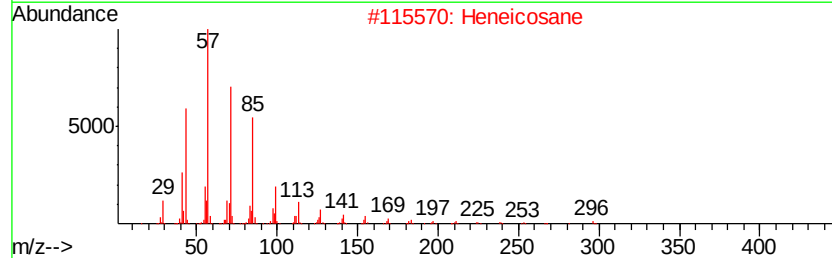
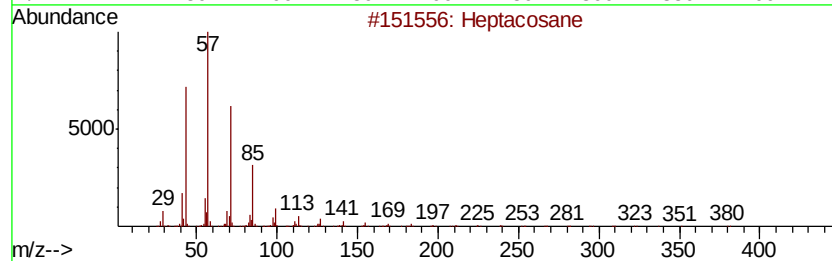
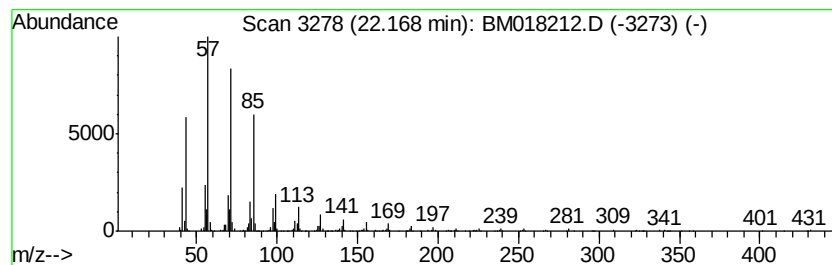
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Heptacosane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.17	9.78 ng	962941	Chrysene-d12	21.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptacosane	380	C27H56	000593-49-7	91
2		Heneicosane	296	C21H44	000629-94-7	91
3		Docosane	310	C22H46	000629-97-0	91
4		Nonacosane	408	C29H60	000630-03-5	91
5		Heptadecane	240	C17H36	000629-78-7	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121518\
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Acq On : 15 Dec 2018 20:08
Operator : JU/SJ
Sample : J6347-06
Misc :
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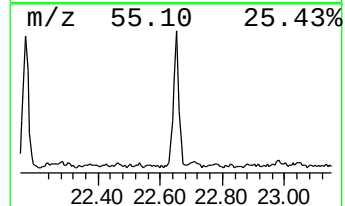
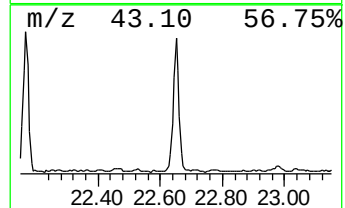
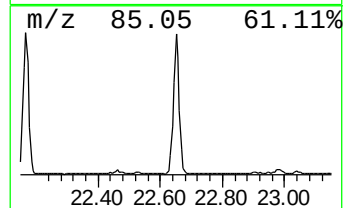
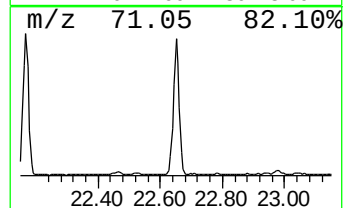
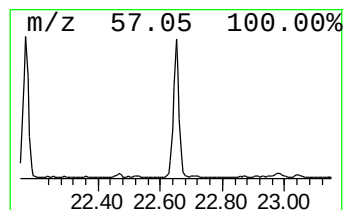
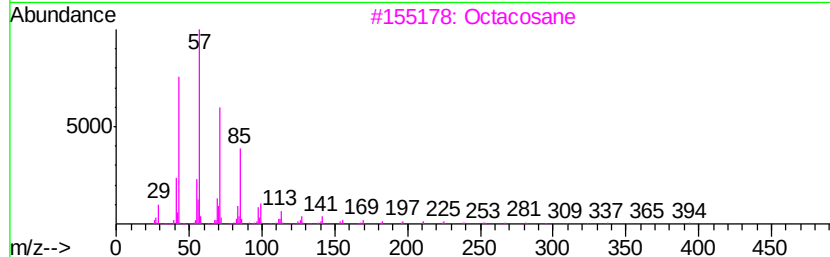
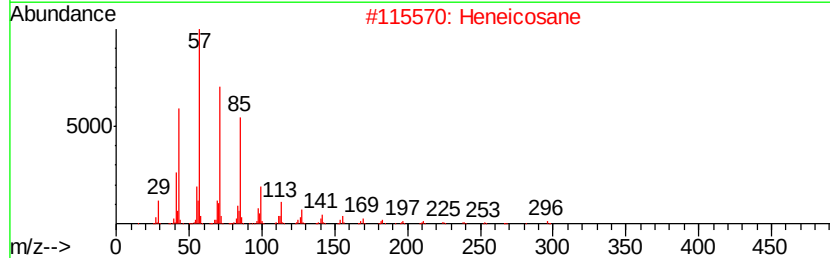
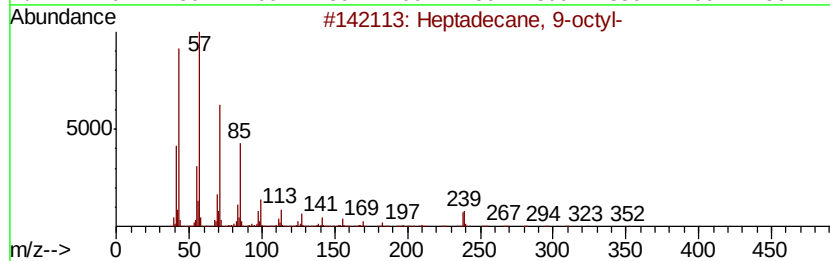
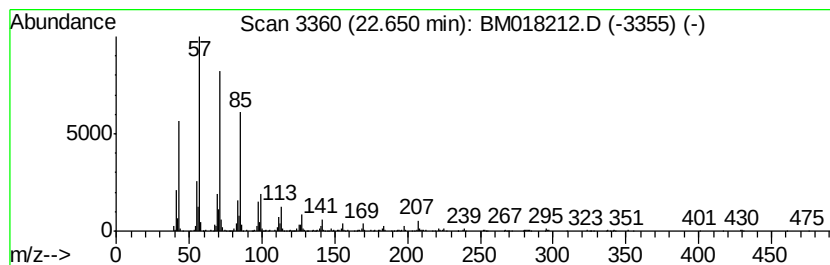
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Heptadecane, 9-octyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.		
22.65	12.87 ng	1038750	Perylene-d12	23.30		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91	
2	Heneicosane	296	C21H44	000629-94-7	91	
3	Octacosane	394	C28H58	000630-02-4	91	
4	Octadecane	254	C18H38	000593-45-3	90	
5	Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	83	



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121518\
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 Acq On : 15 Dec 2018 20:08
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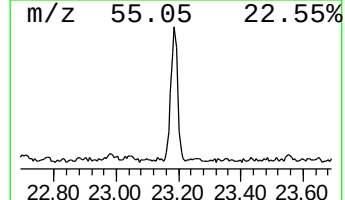
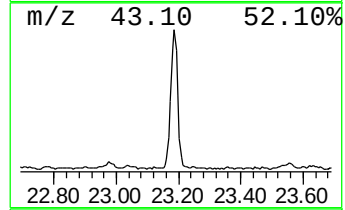
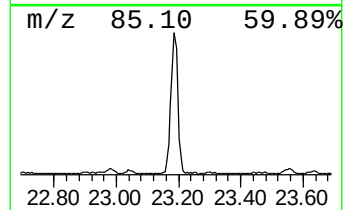
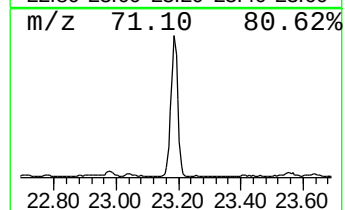
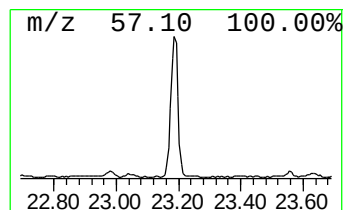
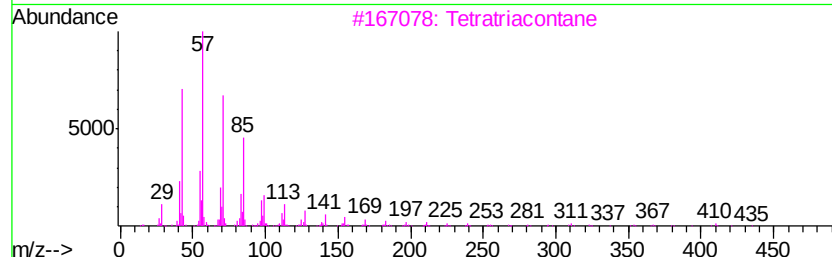
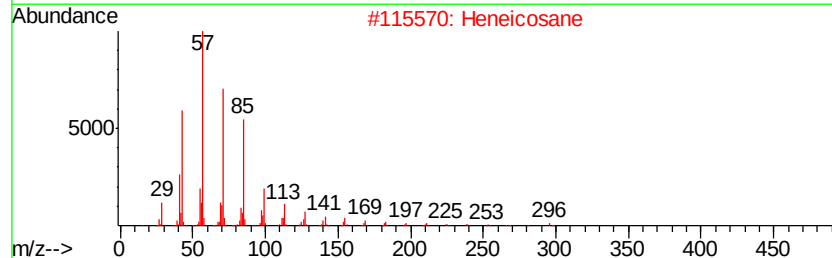
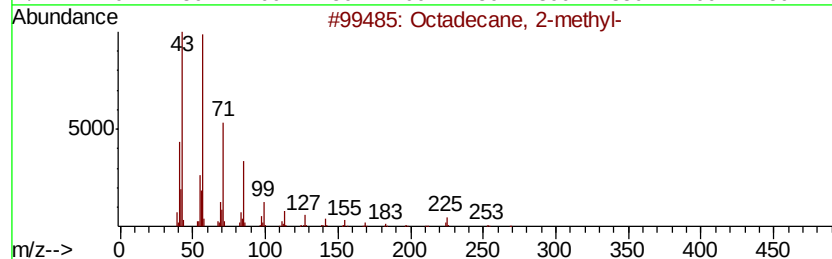
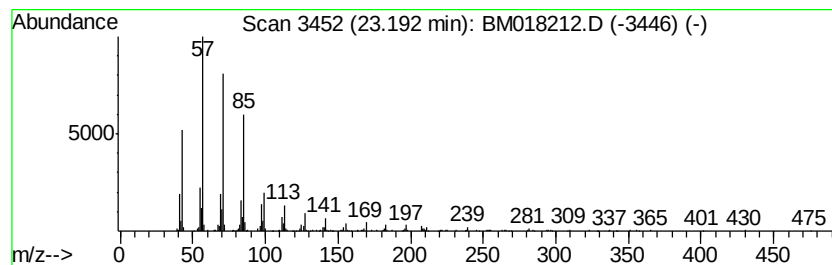
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 11 Octadecane, 2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.19	12.80 ng	1033150	Perylene-d12	23.30

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecane, 2-methyl-	268	C19H40	001560-88-9	91
2		Heneicosane	296	C21H44	000629-94-7	91
3		Tetratriacontane	479	C34H70	014167-59-0	91
4		Heptacosane	380	C27H56	000593-49-7	91
5		Triacotane	422	C30H62	000638-68-6	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121518\
Data File : BM018212.D
Acq On : 15 Dec 2018 20:08
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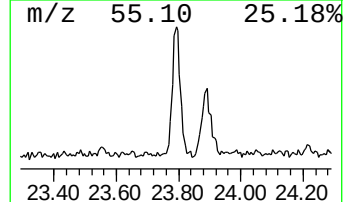
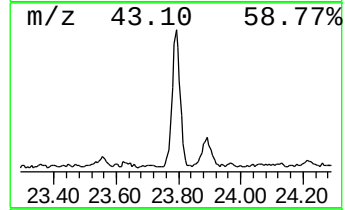
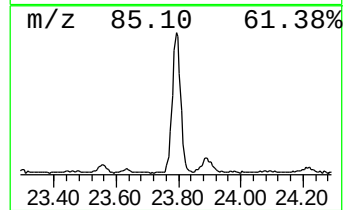
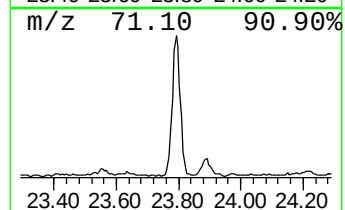
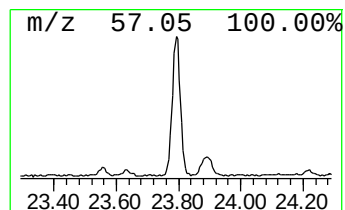
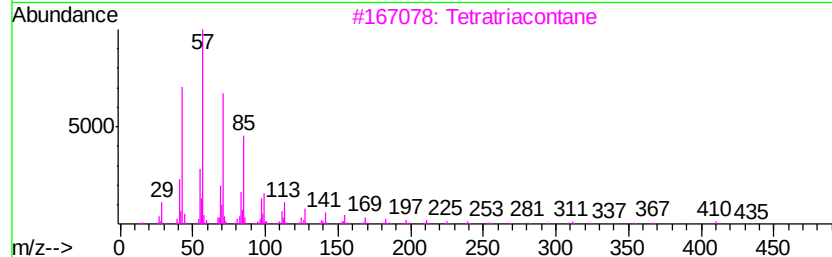
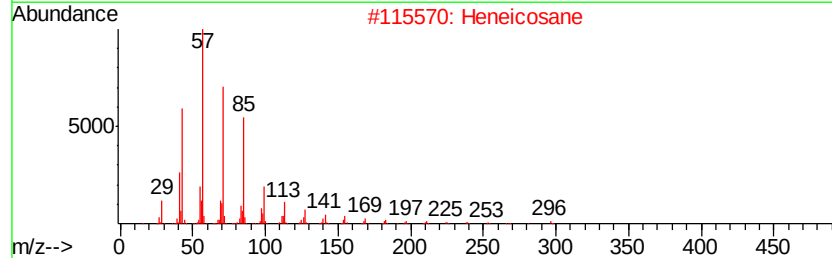
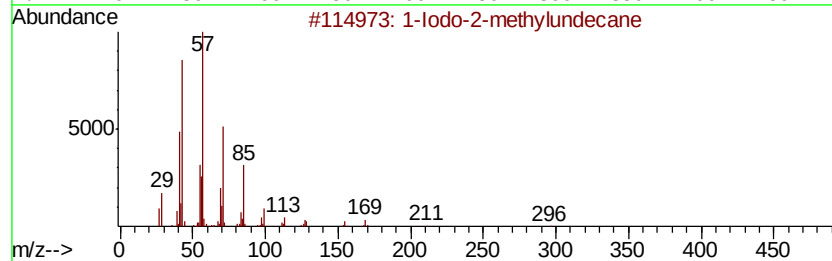
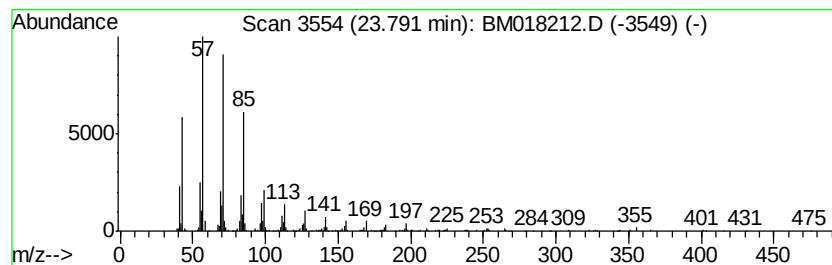
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 1-Iodo-2-methylundecane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.79	8.85 ng	714730	Perylene-d12	23.30

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Iodo-2-methylundecane	296	C12H25I	073105-67-6	94
2			Heneicosane	296	C21H44	000629-94-7	91
3			Tetratriacontane	479	C34H70	014167-59-0	91
4			Triacontane	422	C30H62	000638-68-6	91
5			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121518\
Data File : BM018212.D
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Sample : J6347-06
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\8270-BM121518.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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n-Hexadecanoic acid	17.83	12.6	ng	1055290	4	16.92	1677350	20.0
Octadecanoic acid	19.13	6.0	ng	586580	5	21.14	1969980	20.0
3-Eicosene, (E)-	20.86	10.2	ng	1005910	5	21.14	1969980	20.0
Hexadecane, 3-met...	21.30	4.8	ng	469593	5	21.14	1969980	20.0
Heneicosane	21.72	7.3	ng	723004	5	21.14	1969980	20.0
Heptacosane	22.17	9.8	ng	962941	5	21.14	1969980	20.0
Heptadecane, 9-oc...	22.65	12.9	ng	1038750	6	23.30	1614820	20.0
Octadecane, 2-met...	23.19	12.8	ng	1033150	6	23.30	1614820	20.0
1-Iodo-2-methylun...	23.79	8.8	ng	714730	6	23.30	1614820	20.0