

Data Path : Z:\HPCHEM1\BNA F\DATA\BF071116\
 Data File : BF088906.D
 Acq On : 11 Jul 2016 18:27
 Operator : UM/SJ
 Sample : H3915-03
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 P6-B2-0-11-C

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 3 % of largest Peak

Max Peaks: 100

Peak Location: TOP

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

Peak separation: 5

Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF063016.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	1.678	6	8	13	rVV	88544	127882	7.92%	0.767%
2	1.758	13	15	17	rVV	56508	64853	4.02%	0.389%
3	1.793	17	18	28	rVV	615379	1128968	69.92%	6.775%
4	1.930	28	30	38	rVB2	46394	88396	5.47%	0.530%
5	4.867	285	287	293	rVB	64796	98079	6.07%	0.589%
6	5.301	322	325	328	rBV	1149634	1498577	92.81%	8.993%
7	6.353	414	417	420	rVB	1362976	1367818	84.72%	8.208%
8	6.479	425	428	430	rBV	1695886	1526471	94.54%	9.160%
9	6.707	446	448	450	rVB	425279	398667	24.69%	2.392%
10	6.867	459	462	464	rBV	1470678	1233233	76.38%	7.401%
11	7.279	495	498	500	rBV	1061386	1075821	66.63%	6.456%
12	7.999	558	561	563	rBB	527399	535218	33.15%	3.212%
13	9.073	652	655	657	rBV	1853884	1614602	100.00%	9.689%
14	9.473	687	690	692	rBB	176446	228288	14.14%	1.370%
15	9.748	711	714	716	rBB	523636	526225	32.59%	3.158%
16	10.525	780	782	785	rBV	874392	880984	54.56%	5.287%
17	10.662	790	794	798	rBV	35717	47060	2.91%	0.282%
18	11.108	828	833	834	rBV	37402	33022	2.05%	0.198%
19	11.222	841	843	845	rBV	481992	463503	28.71%	2.781%
20	12.422	946	948	950	rBV	31083	25647	1.59%	0.154%
21	12.685	965	971	979	rBV	574768	915449	56.70%	5.494%
22	12.799	979	981	984	rVB	1035809	1176694	72.88%	7.061%
23	13.702	1058	1060	1064	rVB2	69119	96831	6.00%	0.581%
24	13.839	1069	1072	1078	rBV	501672	508577	31.50%	3.052%
25	14.342	1114	1116	1120	rVB3	15812	32129	1.99%	0.193%
26	14.720	1147	1149	1151	rVB	16324	17736	1.10%	0.106%
27	14.834	1156	1159	1163	rBV	39810	82913	5.14%	0.498%
28	14.925	1165	1167	1171	rBV3	16972	35123	2.18%	0.211%
29	15.097	1180	1182	1185	rVB	58008	79266	4.91%	0.476%
30	15.154	1185	1187	1190	rBV	26862	34877	2.16%	0.209%
31	15.211	1190	1192	1195	rBV	298424	349532	21.65%	2.098%
32	15.543	1218	1221	1224	rVB	54800	96194	5.96%	0.577%
33	15.691	1232	1234	1238	rVB4	11580	22428	1.39%	0.135%
34	16.057	1262	1266	1273	rBV2	52209	112543	6.97%	0.675%

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Instrument :
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ClientSampleId :
P6-B2-0-11-C

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

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

Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF063016.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	16.674	1316	1320	1325	rBV	36377	88561	5.49%	0.531%
36	17.440	1383	1387	1394	rVB2	17240	51691	3.20%	0.310%

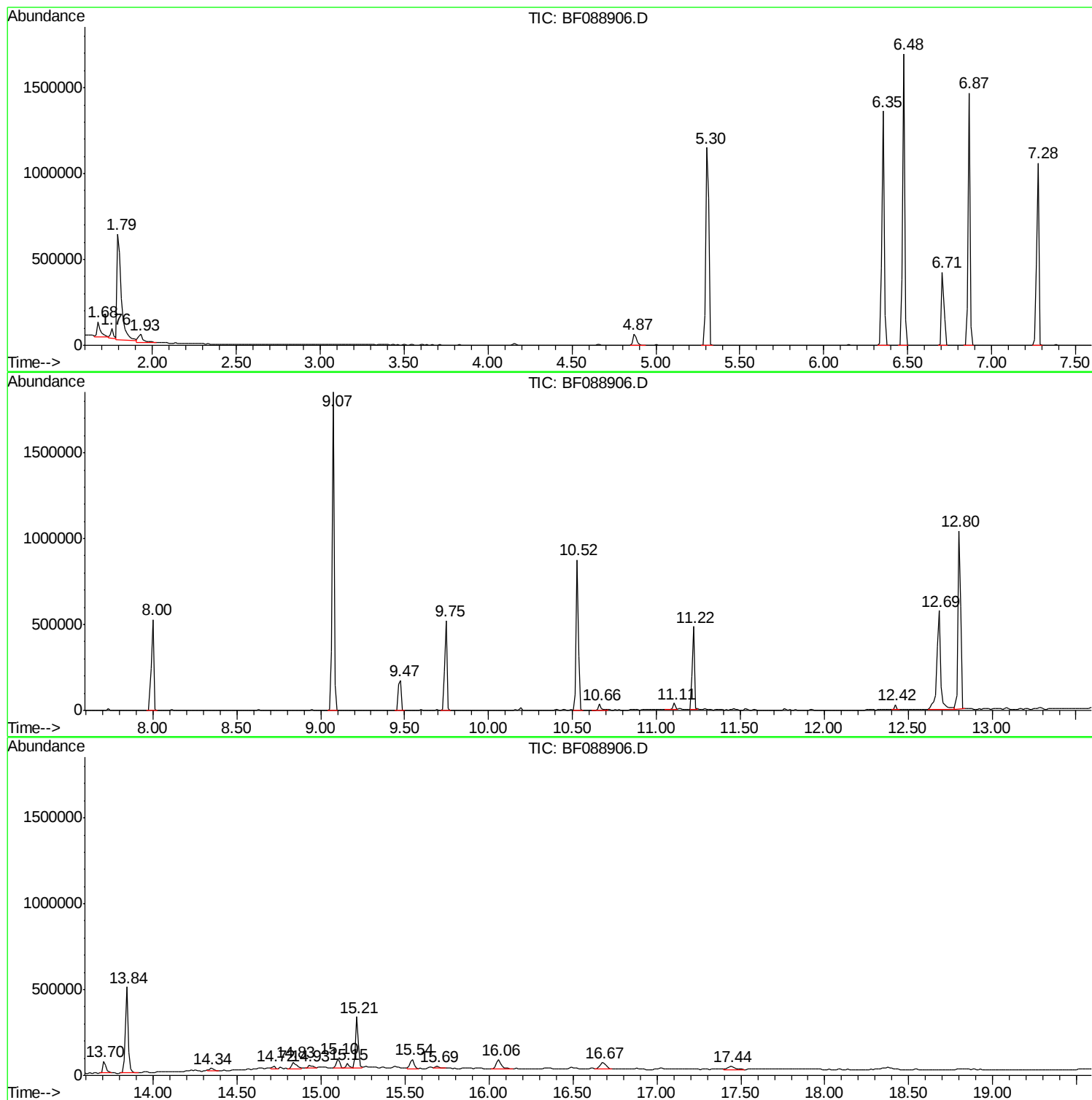
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF063016.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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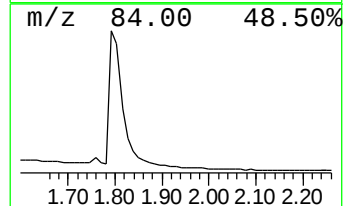
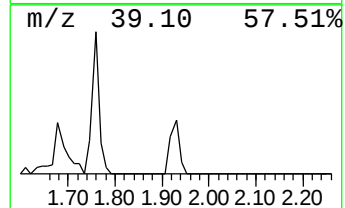
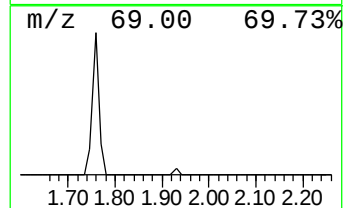
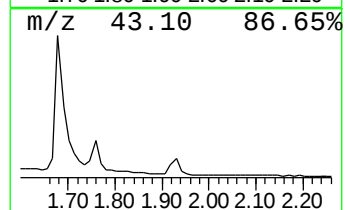
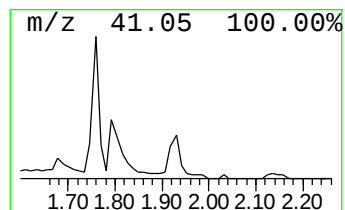
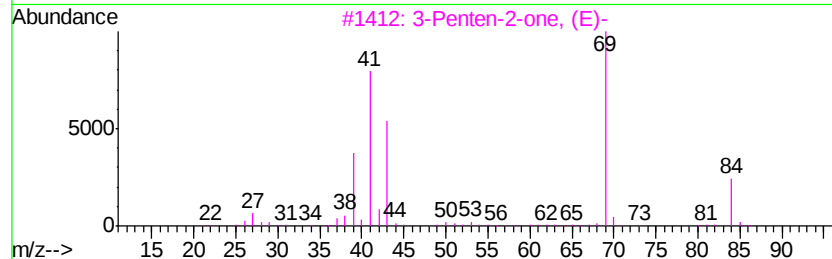
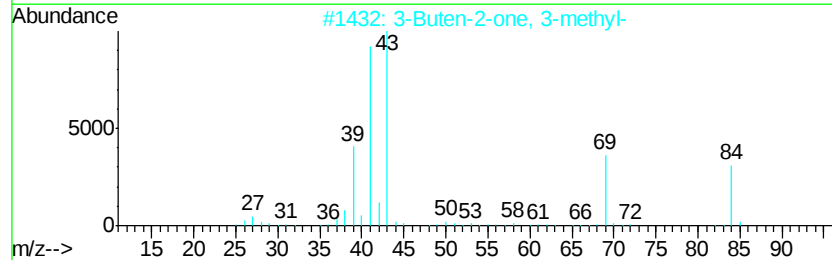
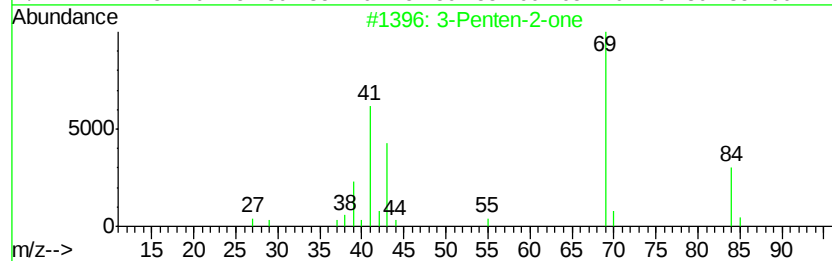
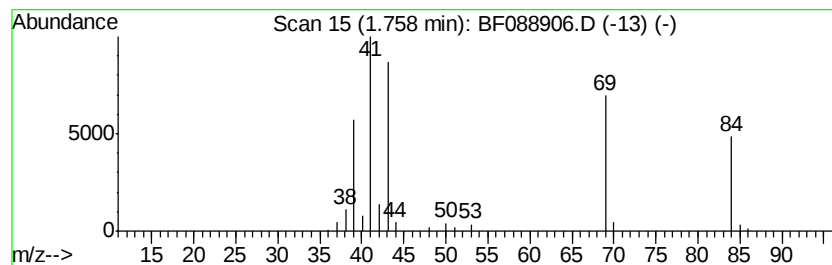
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF063016.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 3-Penten-2-one Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.76	3.25 ng	64853	1,4-Dichlorobenzene-d4	6.71

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Penten-2-one	84	C5H8O	000625-33-2	86
2			3-Buten-2-one, 3-methyl-	84	C5H8O	000814-78-8	72
3			3-Penten-2-one, (E)-	84	C5H8O	003102-33-8	52
4			Ethanone, 1-cyclopropyl-	84	C5H8O	000765-43-5	50
5			2-Butene, 2,3-dimethyl-	84	C6H12	000563-79-1	43



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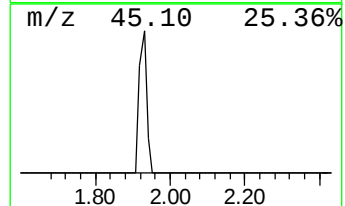
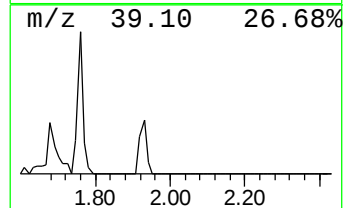
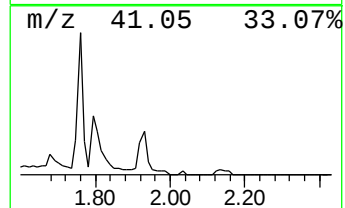
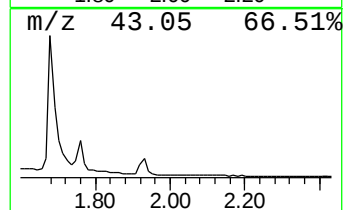
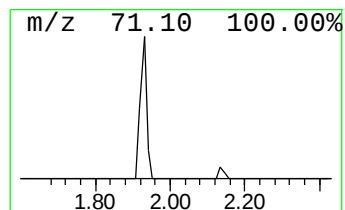
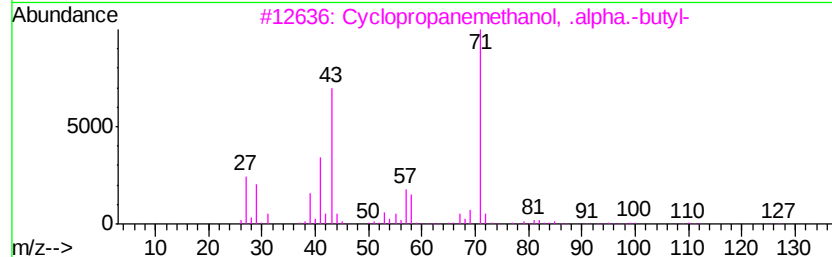
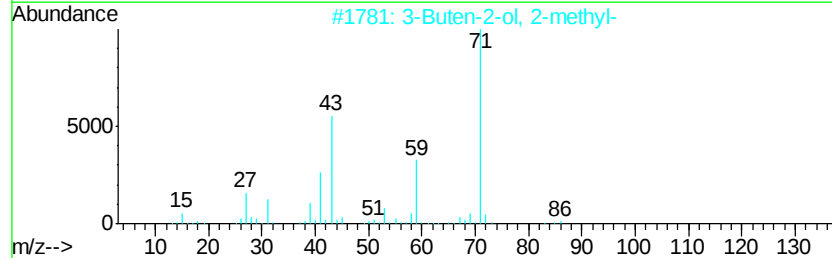
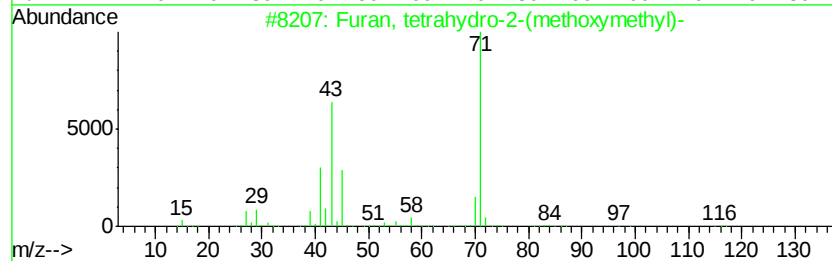
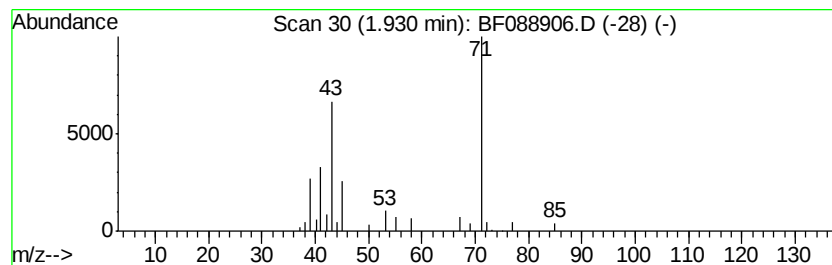
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Furan, tetrahydro-2-(methox... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.93	4.43 ng	88396	1,4-Dichlorobenzene-d4	6.71

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Furan, tetrahydro-2-(methoxymeth...	116	C6H12O2	019354-27-9	64
2		3-Buten-2-ol, 2-methyl-	86	C5H10O	000115-18-4	45
3		Cyclopropanemethanol, .alpha.-bu...	128	C8H16O	004379-16-2	42
4		2,2'-Bifuran, octahydro-	142	C8H14O2	001592-33-2	40
5		2-Hexene, 1-methoxy-, (E)-	114	C7H14O	056052-83-6	39



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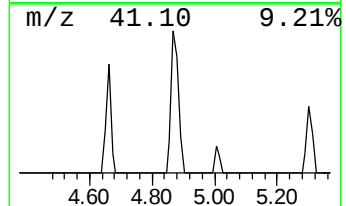
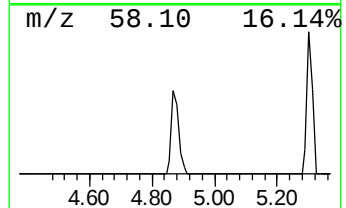
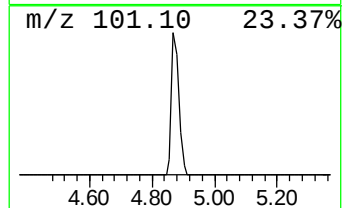
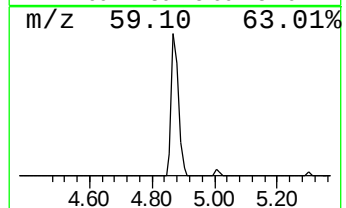
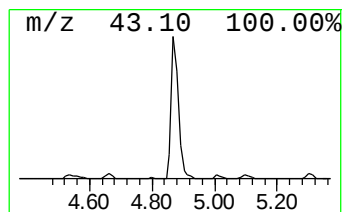
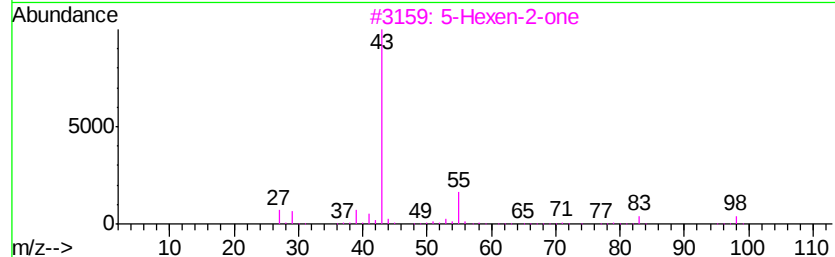
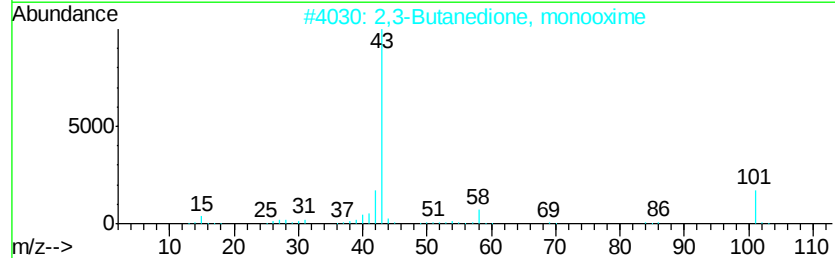
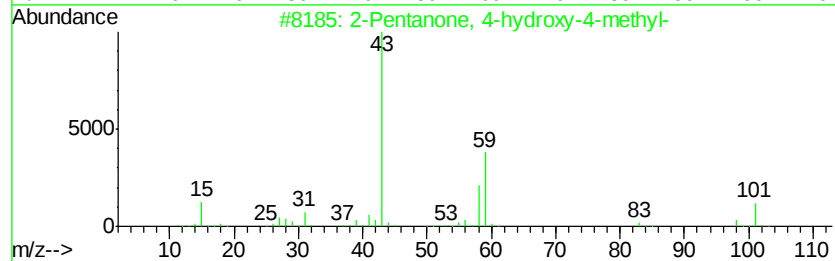
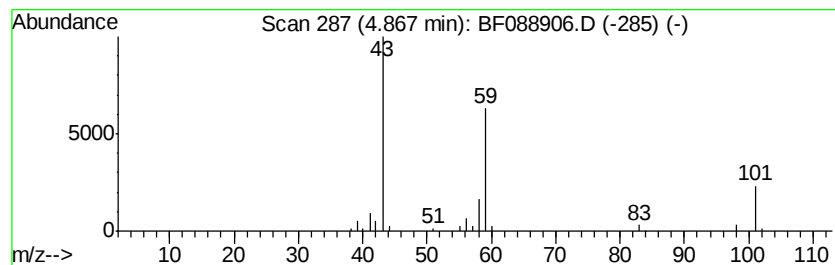
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.87	4.92 ng	98079	1,4-Dichlorobenzene-d4	6.71

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50		
2	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	12		
3	5-Hexen-2-one	98	C6H10O	000109-49-9	9		
4	Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9		
5	(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	9		



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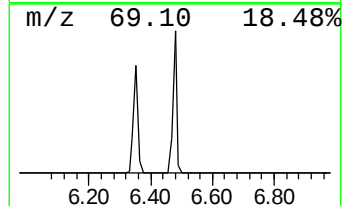
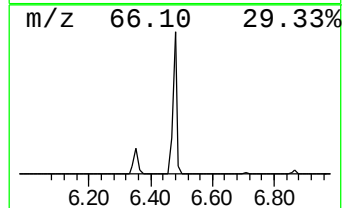
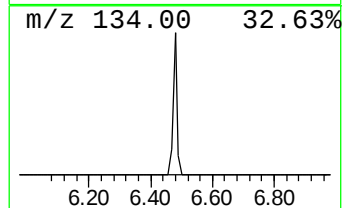
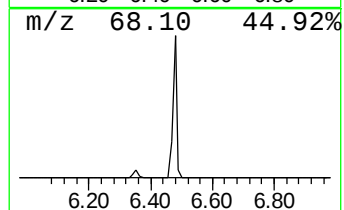
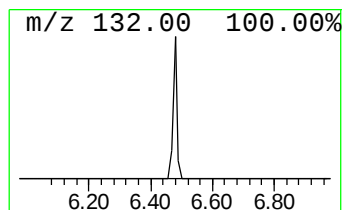
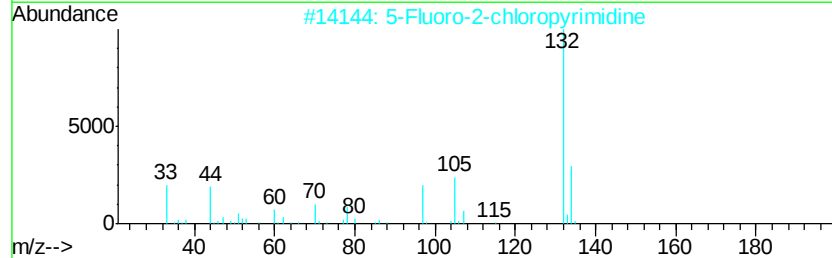
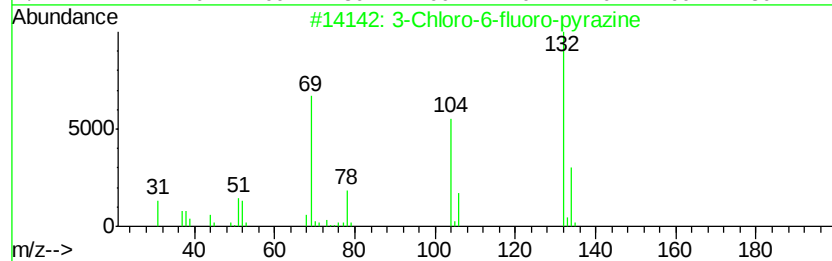
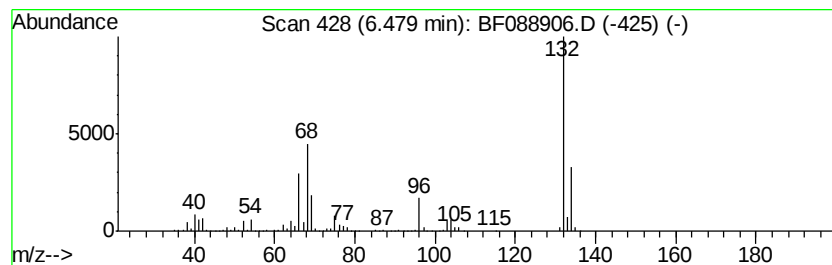
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 unknown6.48 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.48	76.58 ng	1526470	1,4-Dichlorobenzene-d4	6.71

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		Tranvlcyproline-propionyl	189	C12H15NO	1000123-86-3	10
4		N-(-3,4-Methylenedioxyphenylisop...	267	C17H17NO2	1000378-96-6	10
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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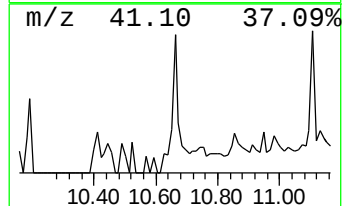
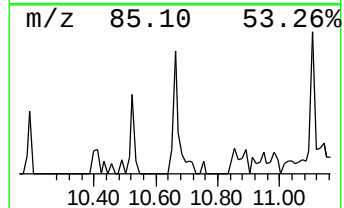
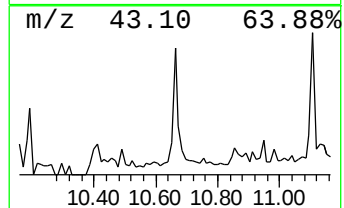
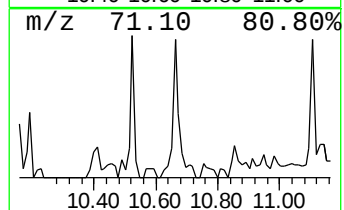
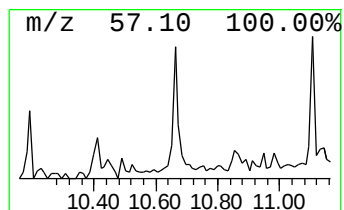
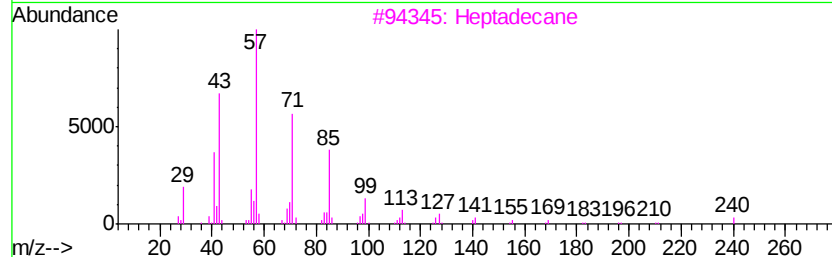
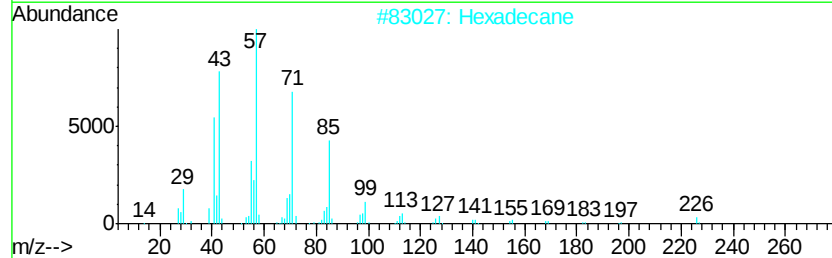
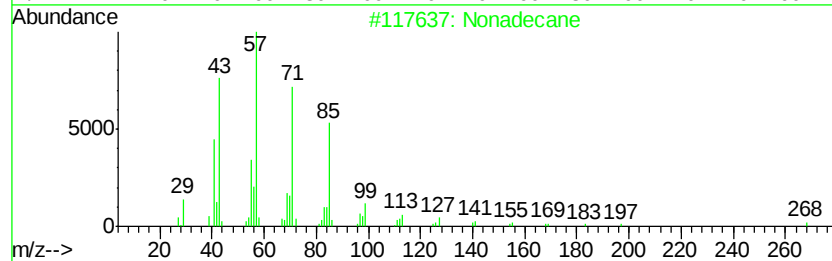
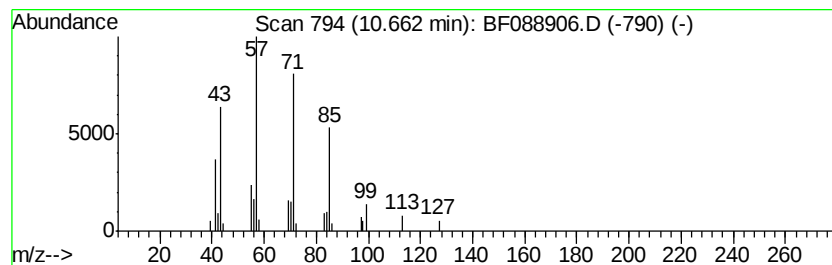
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TIC Library : C:\Database\NIST11.L
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Peak Number 8 Nonadecane Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.66	2.03 ng	47060	Phenanthrene-d10	11.22

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonadecane		268	C19H40	000629-92-5	91
2	Hexadecane		226	C16H34	000544-76-3	90
3	Heptadecane		240	C17H36	000629-78-7	90
4	Heptacosane		380	C27H56	000593-49-7	83
5	Undecane, 5,7-dimethyl-		184	C13H28	017312-83-3	72



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Misc :
ALS Vial : 10 Sample Multiplier: 1

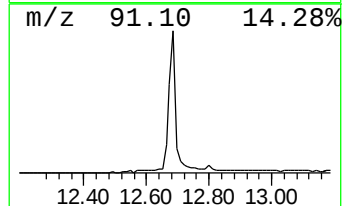
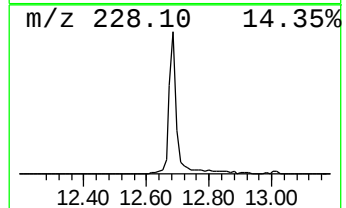
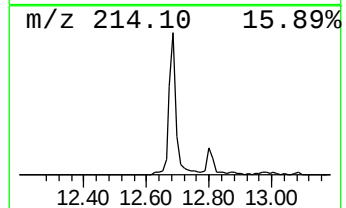
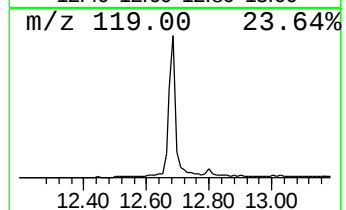
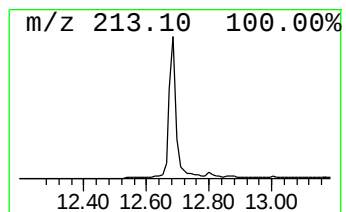
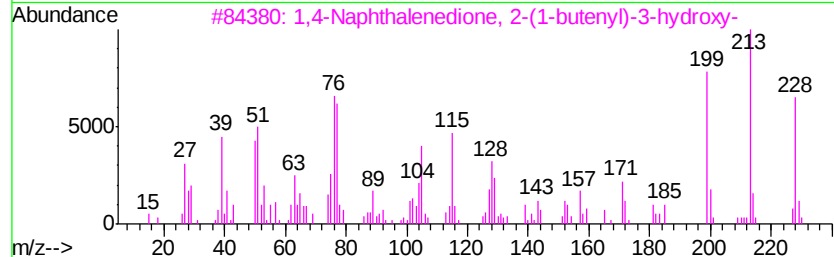
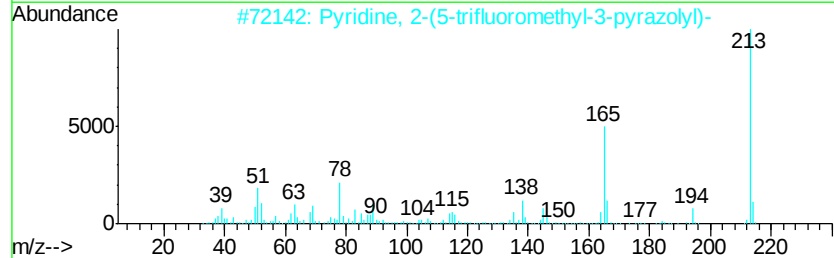
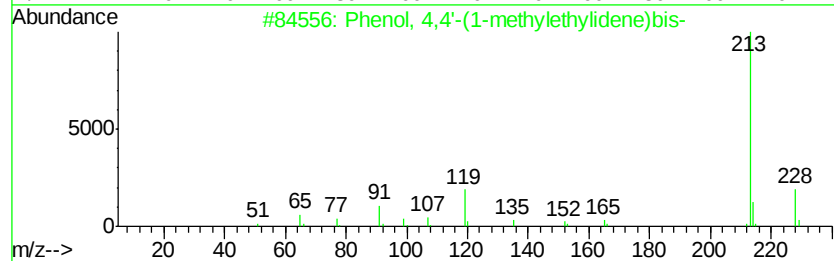
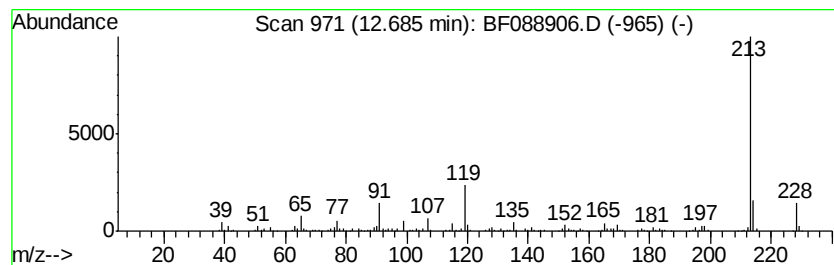
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ClientSampleId :
P6-B2-0-11-C

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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 Phenol, 4,4'-(1-methylethyl... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.69	36.00 ng	915449	Chrysene-d12	13.84	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 4,4'-(1-methylethylidene)bis-	228	C15H16O2	000080-05-7	96
2	Pyridine, 2-(5-trifluoromethyl-3-pyrazolyl)-	213	C9H6F3N3	003974-71-8	50
3	1,4-Naphthalenedione, 2-(1-butenyl)-3-hydroxy-	228	C14H12O3	029366-43-6	47
4	2,4-(1H,3H)-Pyrimidinedione, 1,3-bis-(4-methyl-5-oxo-1,2,4-triazol-2-yl)-	228	C9H16N2O3Si	089447-84-7	42
5	1,4-Puran, 5,6-dihydro-3-[3-indolyl]-	213	C14H15NO	1000128-72-9	40



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF071116\
Data File : BF088906.D
Acq On : 11 Jul 2016 18:27
Operator : UM/SJ
Sample : H3915-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

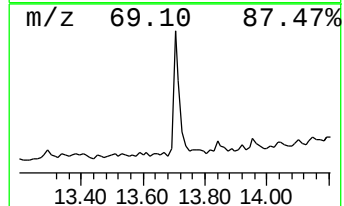
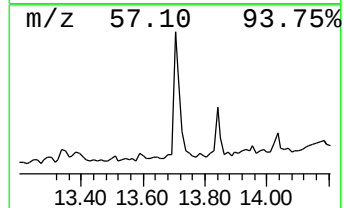
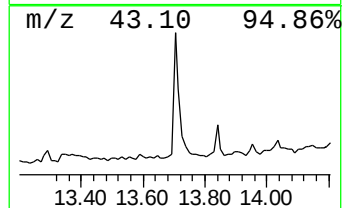
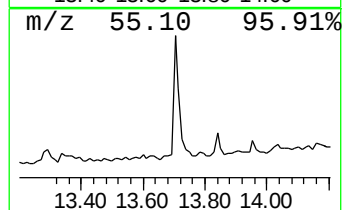
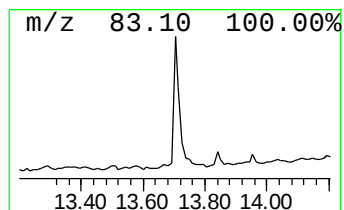
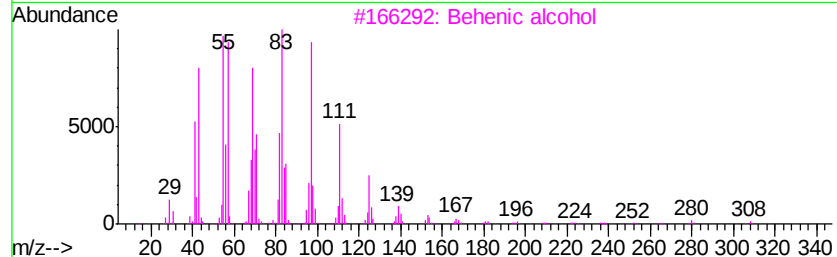
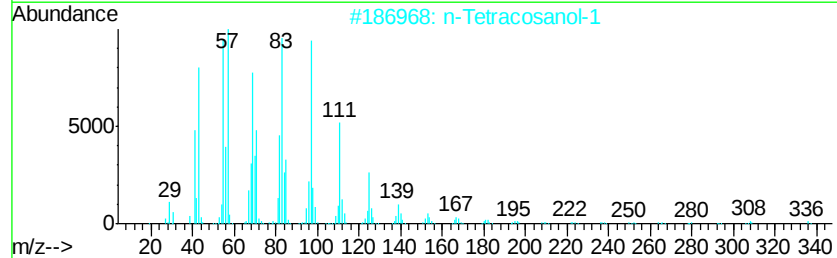
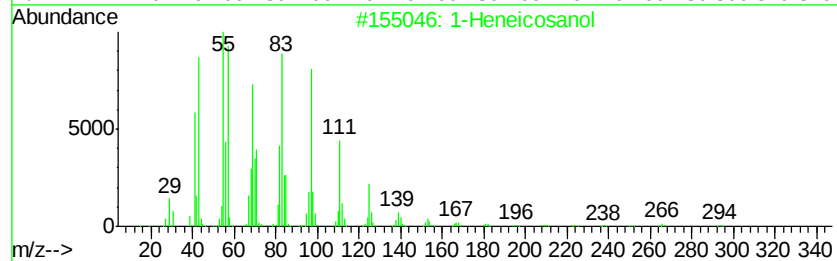
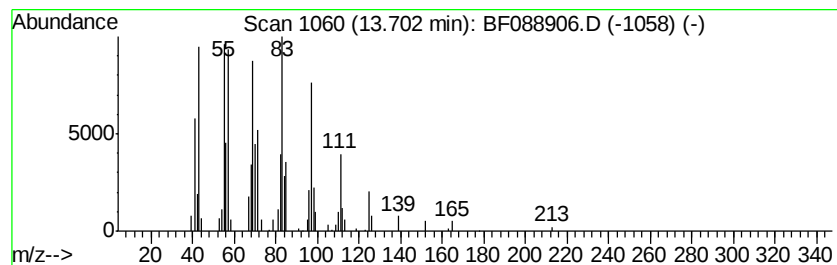
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 1-Heneicosanol Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD		R.T.
13.70	3.81 ng	96831	Chrysene-d12		13.84
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heneicosanol	312	C21H44O	015594-90-8	95
2	n-Tetracosanol-1	354	C24H50O	000506-51-4	94
3	Behenic alcohol	326	C22H46O	000661-19-8	94
4	1-Heptacosanol	396	C27H56O	002004-39-9	91
5	3-Chloropropionic acid, heptadec...	346	C20H39ClO2	1000283-05-1	91



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF071116\
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Acq On : 11 Jul 2016 18:27
Operator : UM/SJ
Sample : H3915-03
Misc :
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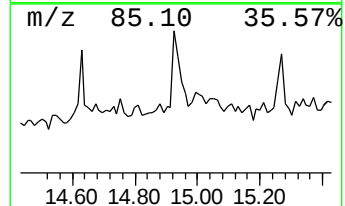
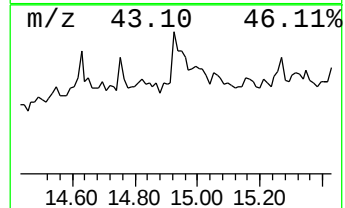
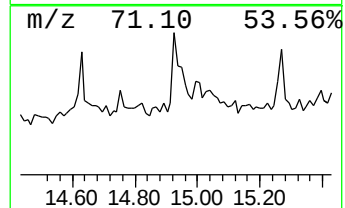
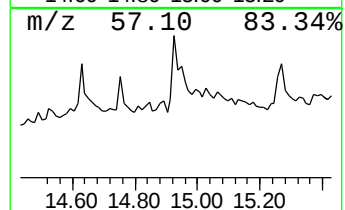
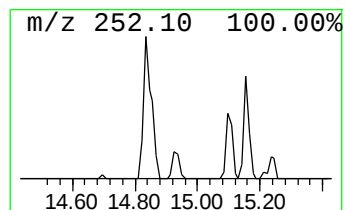
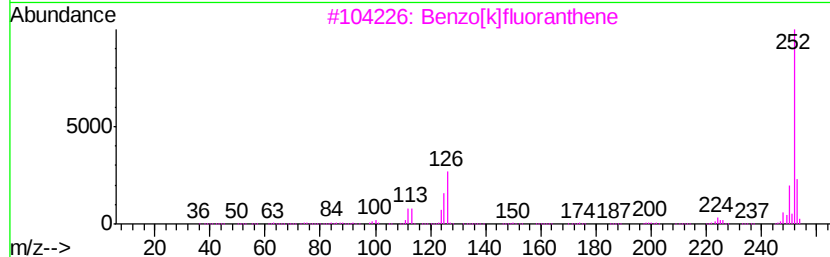
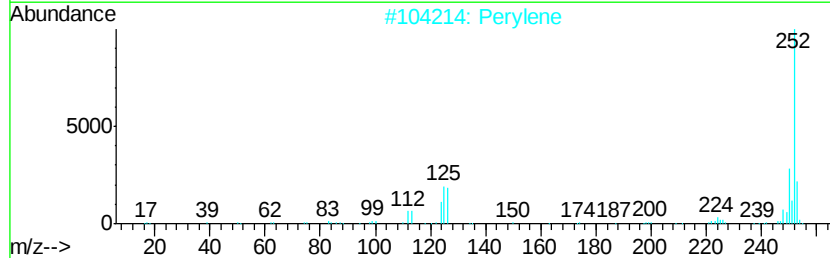
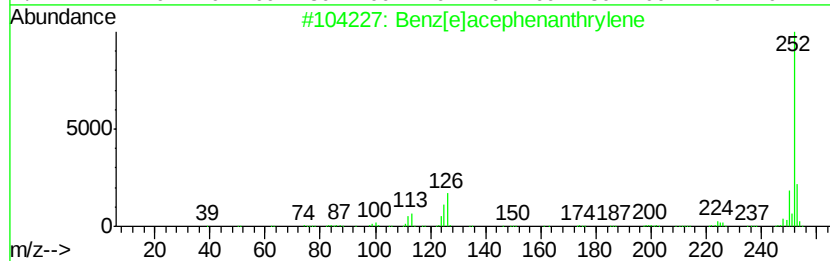
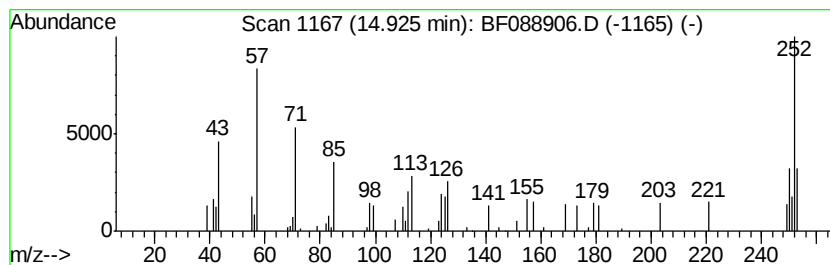
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF063016.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Benz[e]acephenanthrylene Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.93	2.01 ng	35123	Perylene-d12	15.21	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benz[e]acephenanthrylene	252	C20H12	000205-99-2	64
2	Perylene	252	C20H12	000198-55-0	52
3	Benzo[k]fluoranthene	252	C20H12	000207-08-9	47
4	Benzo[i]fluoranthene	252	C20H12	000205-82-3	38
5	Imidazo[1,2-a]pyridine-3-carboxa...	252	C15H12N2O2	1000337-52-4	22



Data Path : Z:\HPCHEM1\BNA F\DATA\BF071116\
Data File : BF088906.D
Acq On : 11 Jul 2016 18:27
Operator : UM/SJ
Sample : H3915-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

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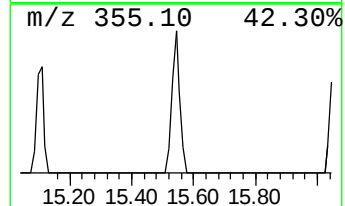
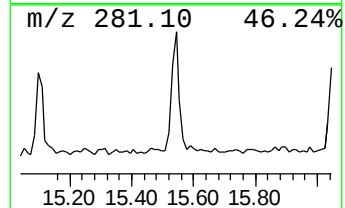
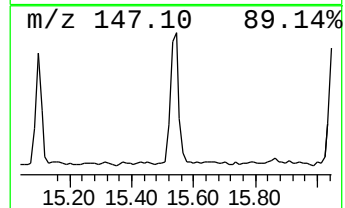
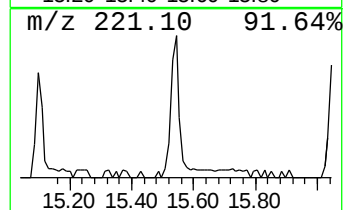
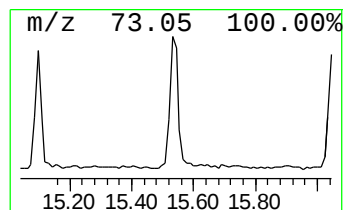
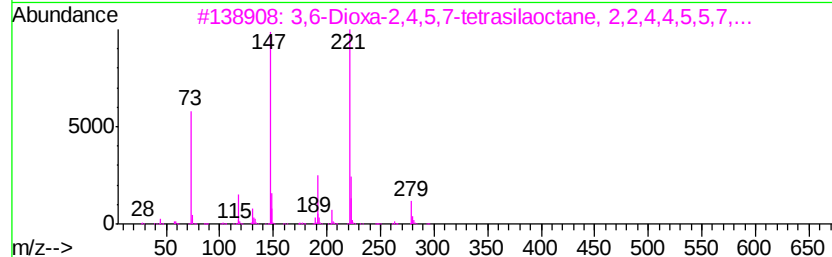
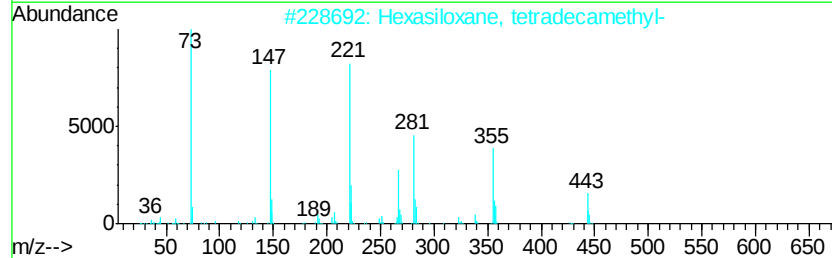
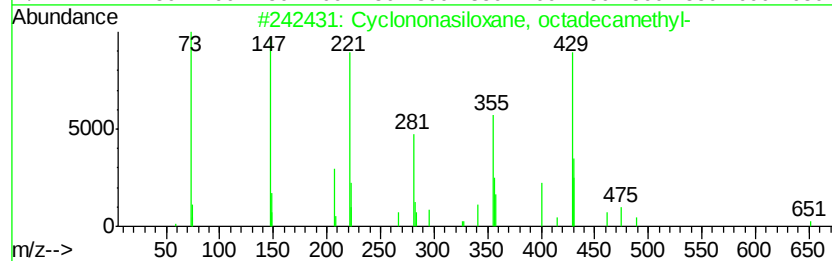
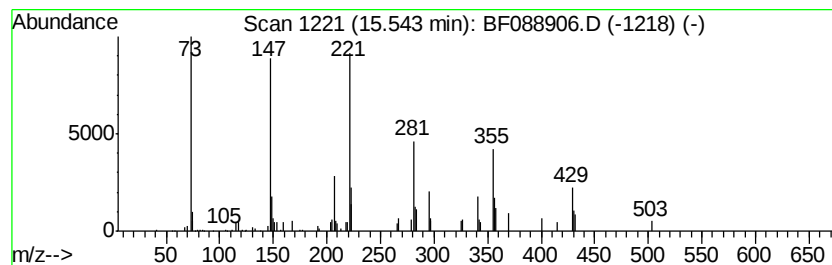
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Cyclononasiloxane, octadeca... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.54	5.50 ng	96194	Perylene-d12	15.21

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclononasiloxane, octadecamethyl-	666	C18H54O9Si9	000556-71-8	87
2			Hexasiloxane, tetradecamethyl-	458	C14H42O5Si6	000107-52-8	81
3			3,6-Dioxa-2,4,5,7-tetrasilaoctan...	294	C10H30O2Si4	004342-25-0	37
4			Mercaptoacetic acid, bis(trimeth...	236	C8H20O2SSi2	006398-62-5	37
5			3-Isopropoxy-1,1,1,7,7,7-hexamet...	576	C18H52O7Si7	071579-69-6	25



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF071116\
Data File : BF088906.D
Acq On : 11 Jul 2016 18:27
Operator : UM/SJ
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Misc :
ALS Vial : 10 Sample Multiplier: 1

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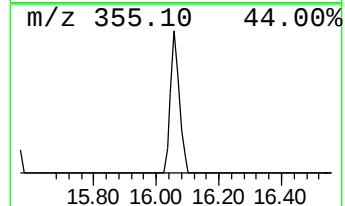
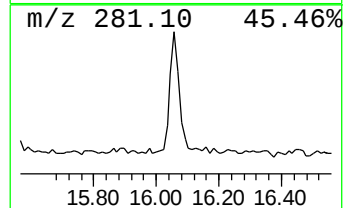
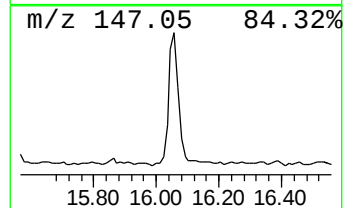
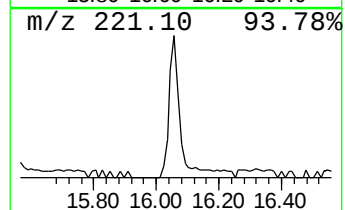
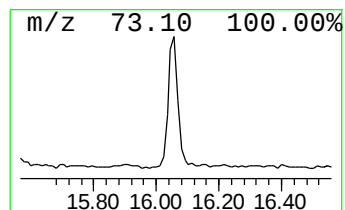
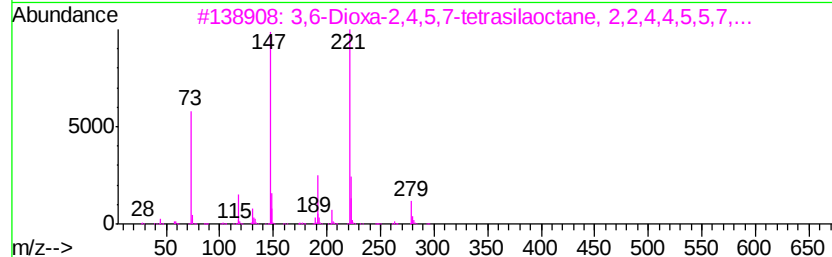
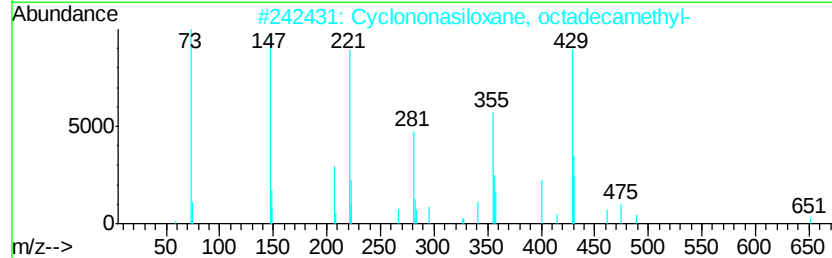
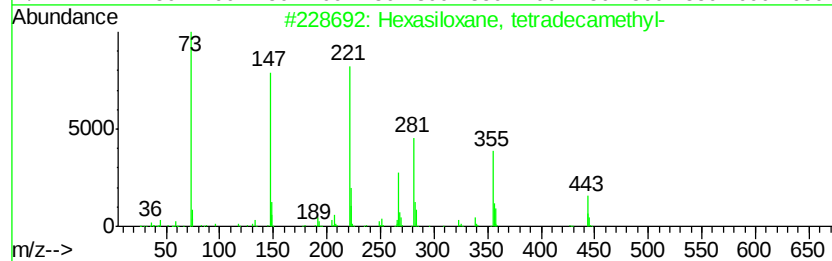
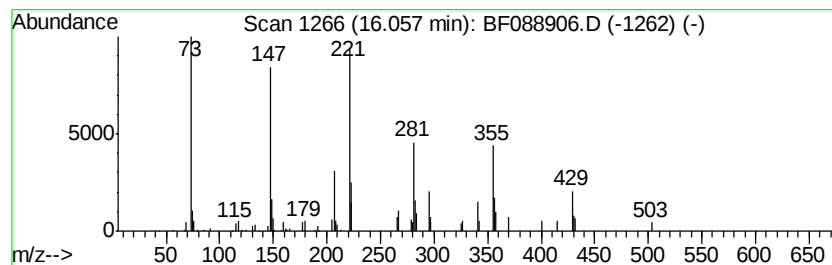
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Hexasiloxane, tetradecamethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.06	6.44 ng	112543	Perylene-d12	15.21

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexasiloxane, tetradecamethyl-	458	C14H42O5Si6	000107-52-8	81
2			Cyclononasiloxane, octadecamethyl-	666	C18H54O9Si9	000556-71-8	64
3			3,6-Dioxa-2,4,5,7-tetrasilaoctan...	294	C10H30O2Si4	004342-25-0	37
4			1,1,1,5,7,7,7-Heptamethyl-3,3-bi...	444	C13H40O5Si6	038147-00-1	28
5			3-Isopropoxy-1,1,1,7,7,7-hexamet...	576	C18H52O7Si7	071579-69-6	25



Data Path : Z:\HPCHEM1\BNA F\DATA\BF071116\
Data File : BF088906.D
Acq On : 11 Jul 2016 18:27
Operator : UM/SJ
Sample : H3915-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

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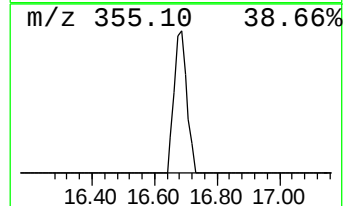
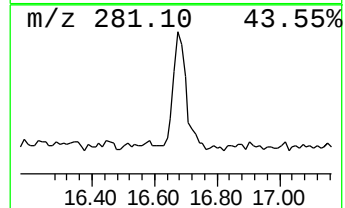
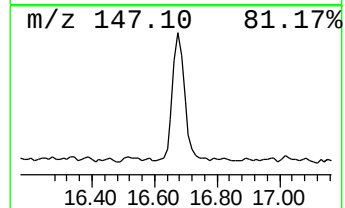
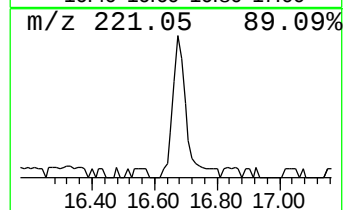
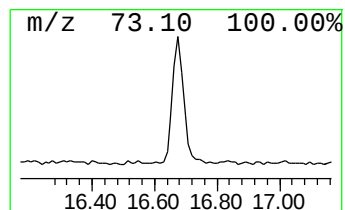
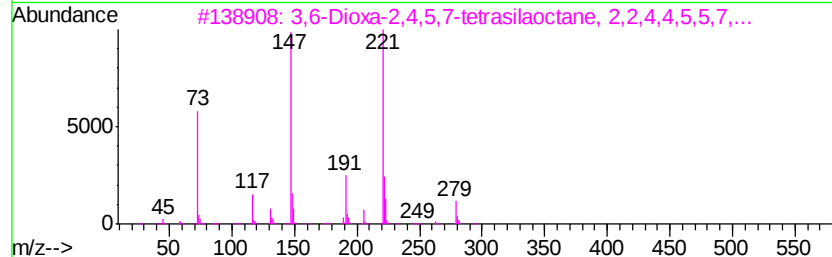
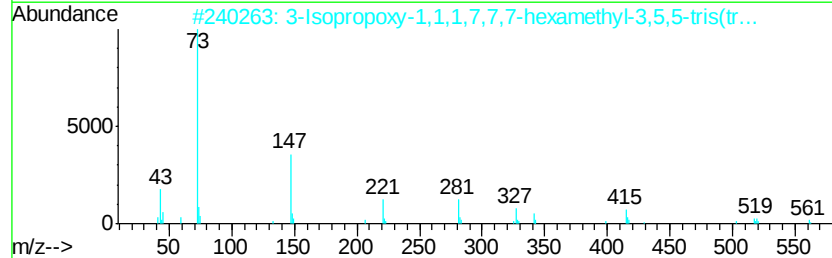
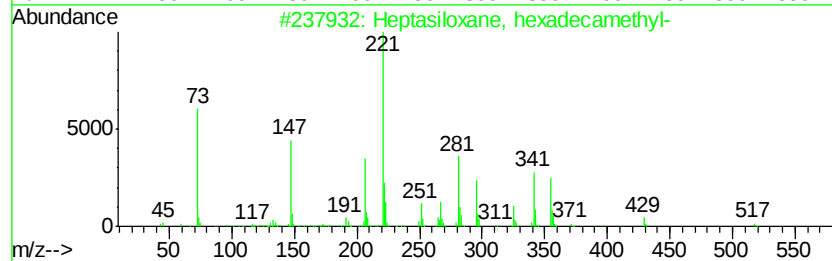
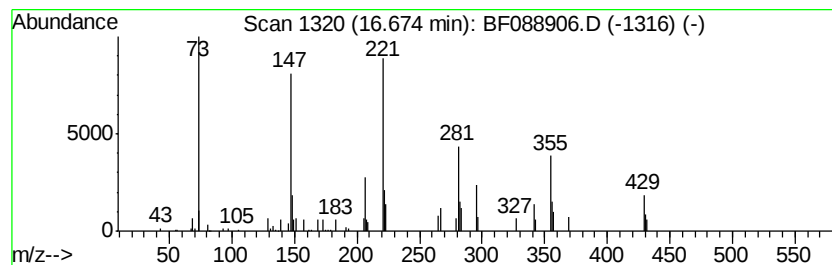
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Heptasiloxane, hexadecamethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.67	5.07 ng	88561	Perylene-d12	15.21

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptasiloxane, hexadecamethyl-	532	C16H48O6Si7	000541-01-5	64
2		3-Isopropoxy-1,1,1,7,7,7-hexamet...	576	C18H52O7Si7	071579-69-6	25
3		3,6-Dioxa-2,4,5,7-tetrasilaoctan...	294	C10H30O2Si4	004342-25-0	25
4		Trisiloxane, 1,1,1,5,5,5-hexamet...	384	C12H36O4Si5	003555-47-3	22
5		N-Benzyl-N-ethyl-p-isopropylbenz...	281	C19H23NO	015089-22-2	22



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF071116\
Data File : BF088906.D
Acq On : 11 Jul 2016 18:27
Operator : UM/SJ
Sample : H3915-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
3-Penten-2-one	1.76	3.3	ng	64853	1	6.71	398667	20.0
Furan, tetrahydro...	1.93	4.4	ng	88396	1	6.71	398667	20.0
2-Pentanone, 4-hy...	4.87	4.9	ng	98079	1	6.71	398667	20.0
unknown6.48	6.48	76.6	ng	1526470	1	6.71	398667	20.0
Nonadecane	10.66	2.0	ng	47060	4	11.22	463503	20.0
Phenol, 4,4'-(1-m...	12.69	36.0	ng	915449	5	13.84	508577	20.0
1-Heneicosanol	13.70	3.8	ng	96831	5	13.84	508577	20.0
Benz[e]acephenant...	14.93	2.0	ng	35123	6	15.21	349532	20.0
Cyclononasiloxane...	15.54	5.5	ng	96194	6	15.21	349532	20.0
Hexasiloxane, tet...	16.06	6.4	ng	112543	6	15.21	349532	20.0
Heptasiloxane, he...	16.67	5.1	ng	88561	6	15.21	349532	20.0