

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032216\
 Data File : BF085671.D
 Acq On : 22 Mar 2016 21:08
 Operator : UM/SJ
 Sample : H1857-02
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-01(23-25)

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-BF031816.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	5.018	236	239	251	rVB	755408	1164567	27.32%	2.932%
2	5.441	272	276	278	rBV	2889775	3645662	85.51%	9.178%
3	5.841	308	311	313	rBV	80592	109668	2.57%	0.276%
4	6.470	363	366	369	rBV	2721610	3226037	75.67%	8.122%
5	6.596	375	377	380	rBV	2537827	3744792	87.84%	9.428%
6	6.836	396	398	400	rBV	594873	741167	17.38%	1.866%
7	6.984	409	411	414	rBV	2345285	2779548	65.20%	6.998%
8	7.396	444	447	450	rBV	1989161	2353576	55.21%	5.925%
9	8.116	507	510	512	rBV	1200842	1054676	24.74%	2.655%
10	9.190	601	604	607	rBV	3064393	4263287	100.00%	10.733%
11	9.590	636	639	641	rBV	607629	585737	13.74%	1.475%
12	9.808	655	658	660	rBV	68328	73584	1.73%	0.185%
13	9.865	661	663	666	rVB	1062427	1148727	26.94%	2.892%
14	10.299	699	701	703	rVB	120498	123844	2.90%	0.312%
15	10.516	717	720	724	rBV	45870	87214	2.05%	0.220%
16	10.665	729	733	735	rBV	2435125	3134988	73.53%	7.893%
17	10.779	741	743	749	rVB	102696	188537	4.42%	0.475%
18	11.225	780	782	783	rBV	52068	64207	1.51%	0.162%
19	11.351	791	793	795	rBV	1474994	1278135	29.98%	3.218%
20	11.888	837	840	848	rBV	207020	329265	7.72%	0.829%
21	12.665	905	908	914	rBV	59624	76852	1.80%	0.193%
22	12.939	929	932	934	rBV	4068449	4193458	98.36%	10.557%
23	13.339	964	967	969	rBV	69324	86296	2.02%	0.217%
24	13.408	969	973	975	rVV3	160056	290070	6.80%	0.730%
25	13.476	975	979	981	rVV	189575	385062	9.03%	0.969%
26	13.511	981	982	985	rVV	49946	60438	1.42%	0.152%
27	13.568	985	987	991	rVB	97300	202843	4.76%	0.511%
28	13.831	1007	1010	1021	rBV	230909	377727	8.86%	0.951%
29	13.979	1021	1023	1026	rVB	1134114	1131305	26.54%	2.848%
30	14.825	1095	1097	1099	rBV	39316	50032	1.17%	0.126%
31	15.397	1144	1147	1150	rBV	608109	810064	19.00%	2.039%
32	16.265	1218	1223	1225	rBV5	376714	1004998	23.57%	2.530%
33	16.311	1225	1227	1228	rVV	278557	371358	8.71%	0.935%
34	16.334	1228	1229	1231	rVB	318653	258463	6.06%	0.651%

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ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-01(23-25)

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-BF031816.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	17.214	1304	1306	1310	rBV3	59325	64606	1.52%	0.163%
36	17.340	1314	1317	1320	rVB5	86680	204356	4.79%	0.514%
37	17.660	1342	1345	1346	rBV2	60836	55330	1.30%	0.139%

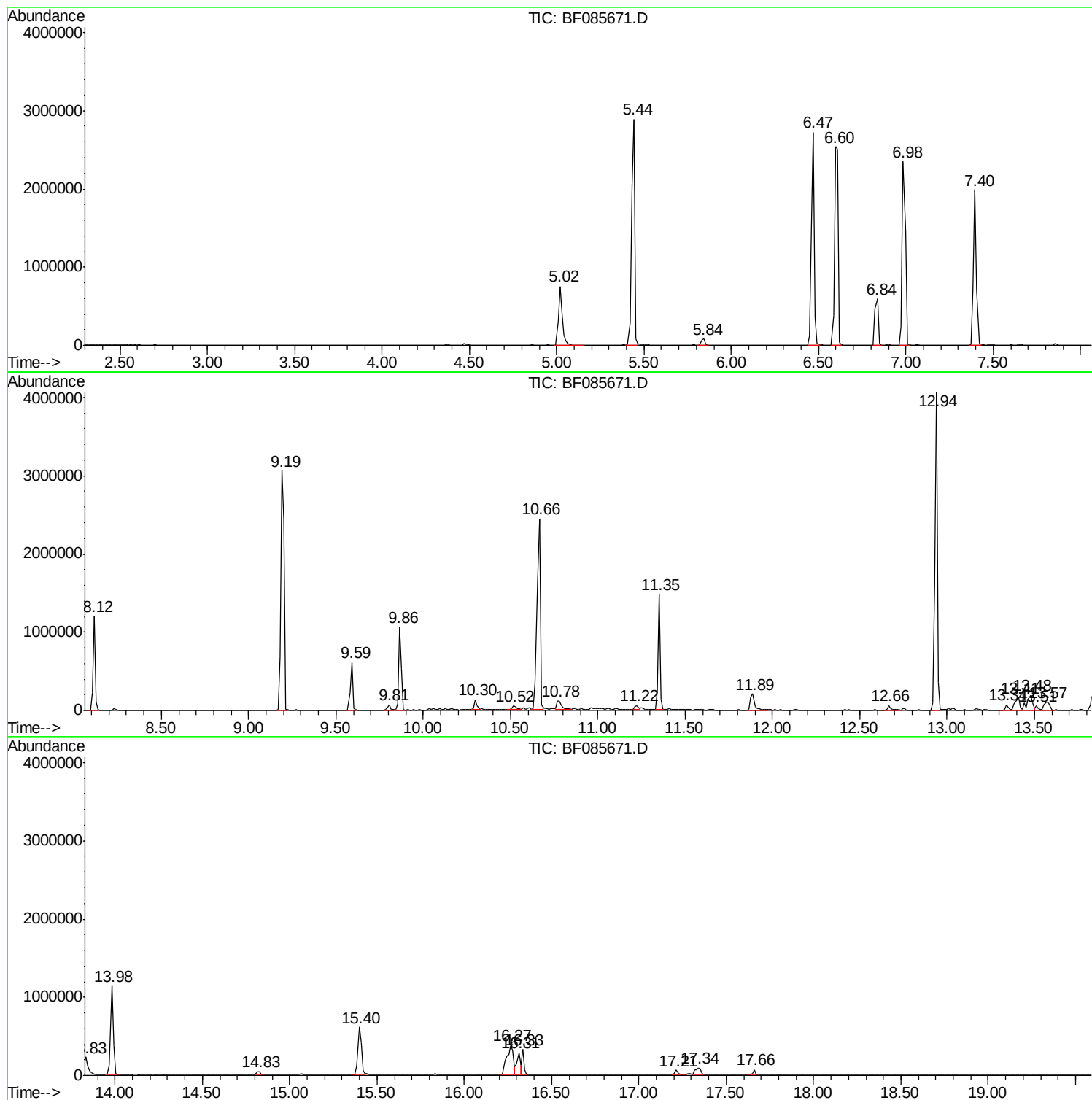
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031816.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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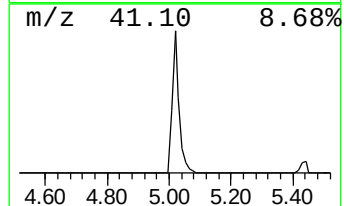
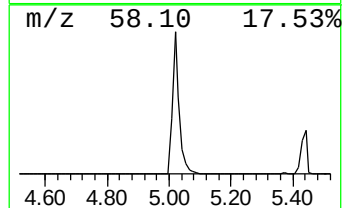
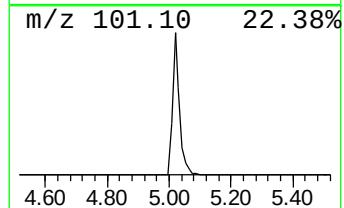
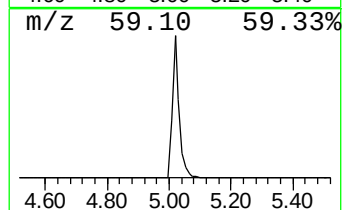
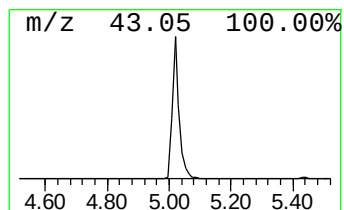
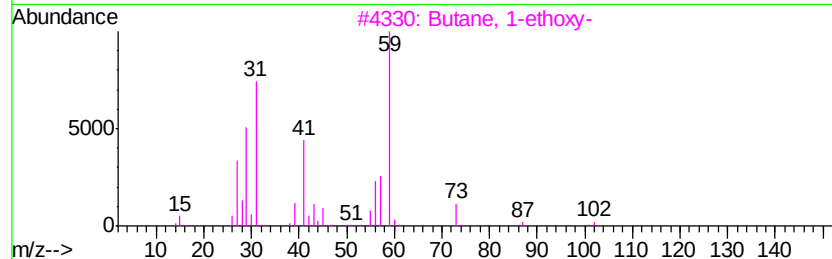
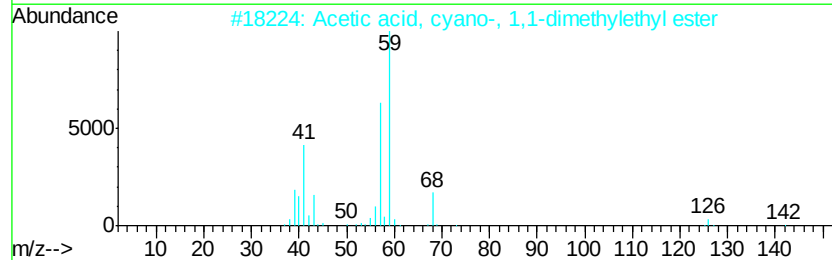
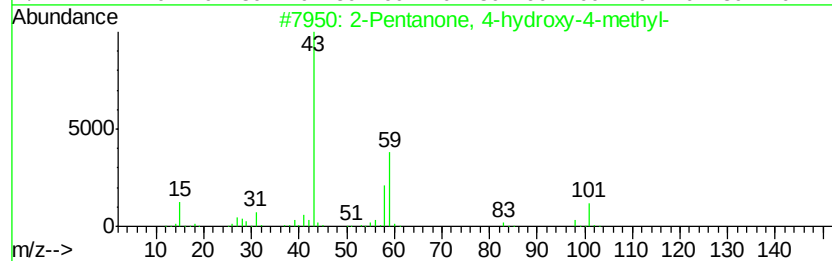
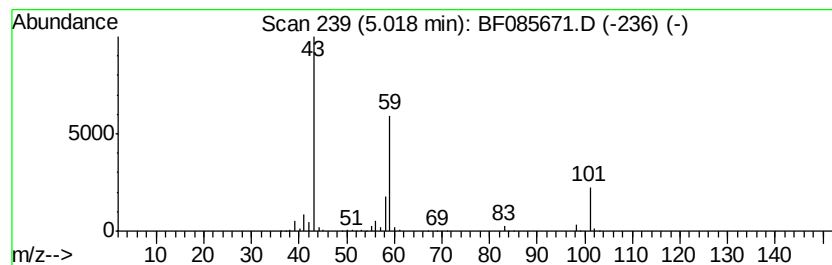
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.02	31.43 ng	1164570	1,4-Dichlorobenzene-d4	6.84

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50		
2	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	17		
3	Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9		
4	5-Hexen-2-one	98	C6H10O	000109-49-9	9		
5	Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9		



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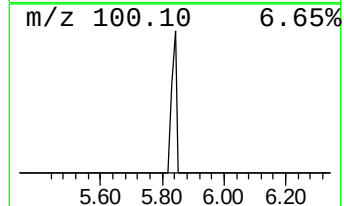
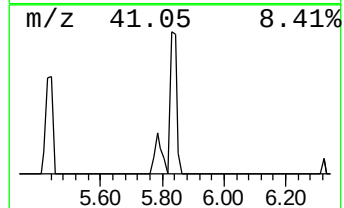
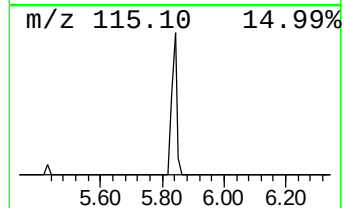
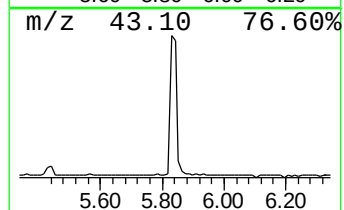
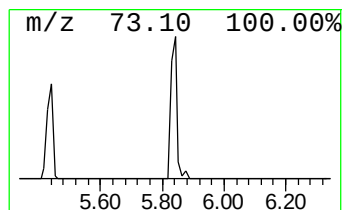
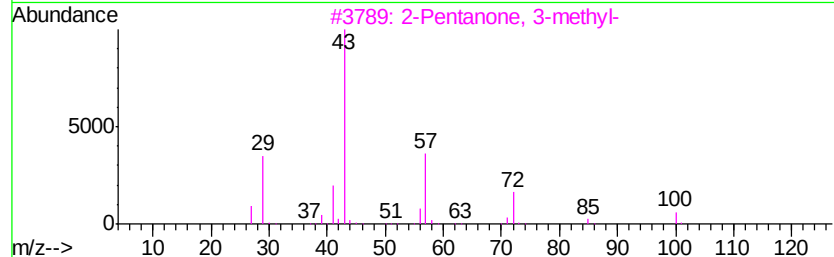
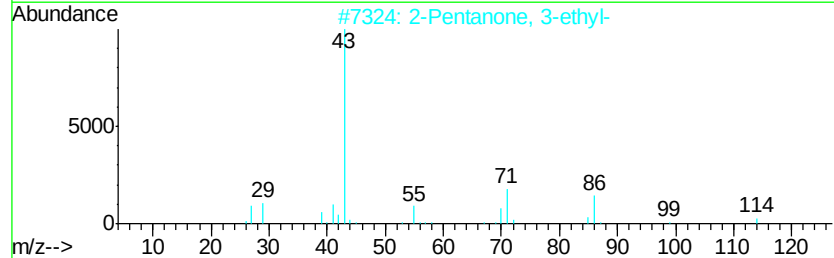
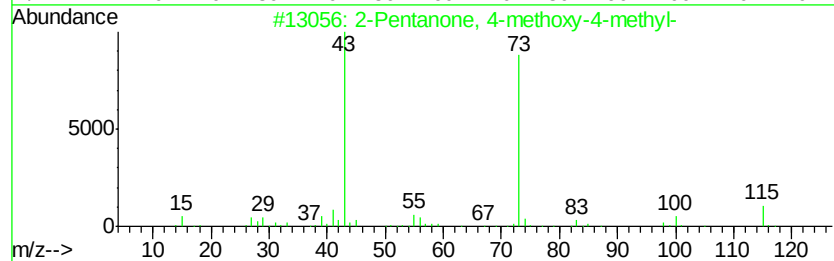
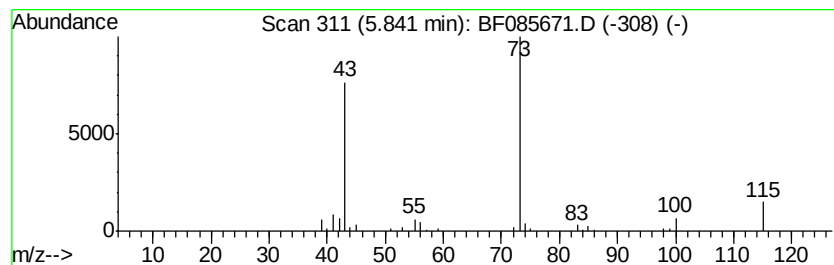
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 2-Pentanone, 4-methoxy-4-me... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.84	2.96 ng	109668	1,4-Dichlorobenzene-d4	6.84

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-methoxy-4-methyl-	130	C7H14O2	000107-70-0	78
2			2-Pentanone, 3-ethyl-	114	C7H14O	006137-03-7	10
3			2-Pentanone, 3-methyl-	100	C6H12O	000565-61-7	9
4			2-Propyn-1-ol, acetate	98	C5H6O2	000627-09-8	9
5			Methyl Isobutyl Ketone	100	C6H12O	000108-10-1	9



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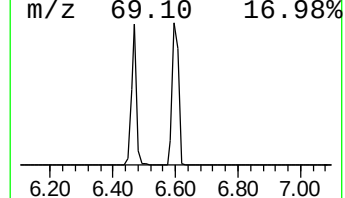
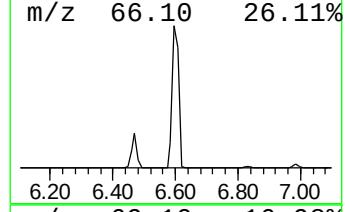
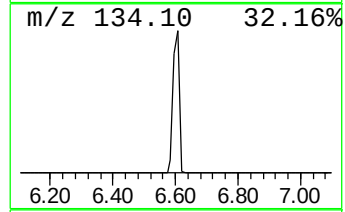
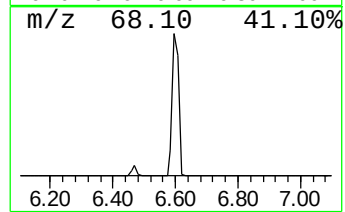
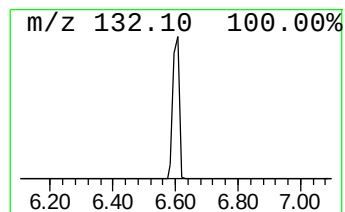
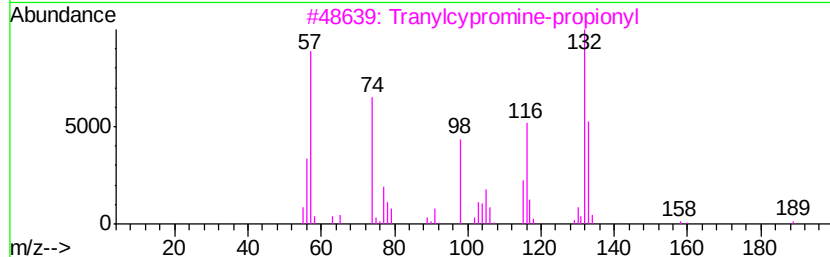
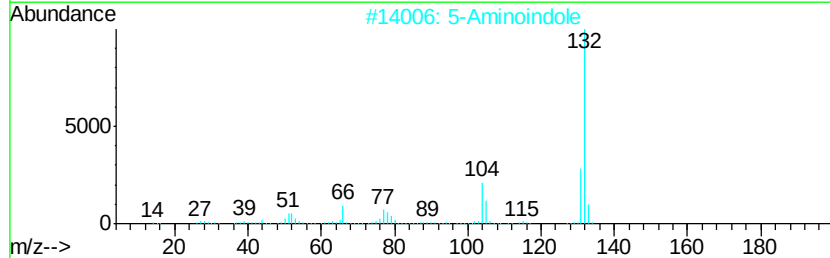
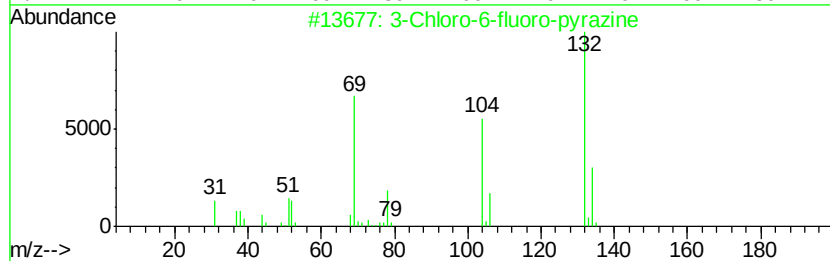
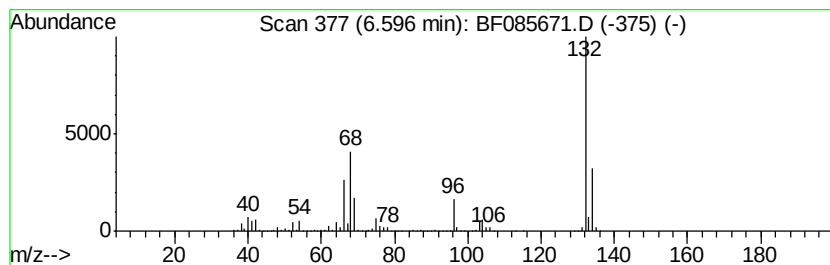
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown6.60 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.60	101.05 ng	3744790	1,4-Dichlorobenzene-d4	6.84

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		5-Aminoindole	132	C8H8N2	005192-03-0	12
3		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
4		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9
5		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9



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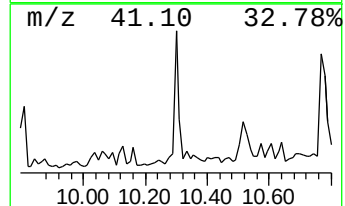
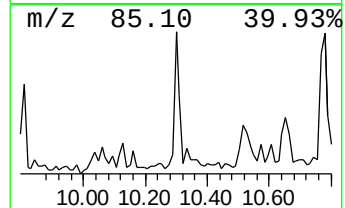
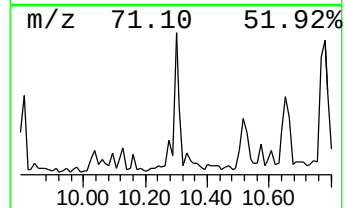
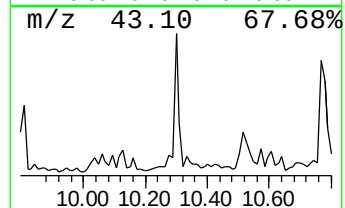
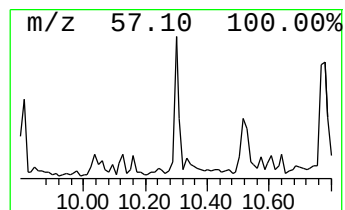
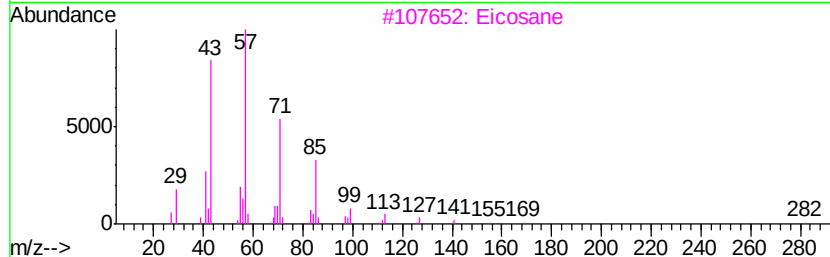
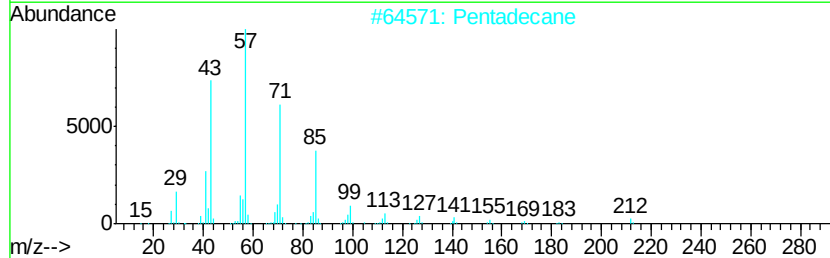
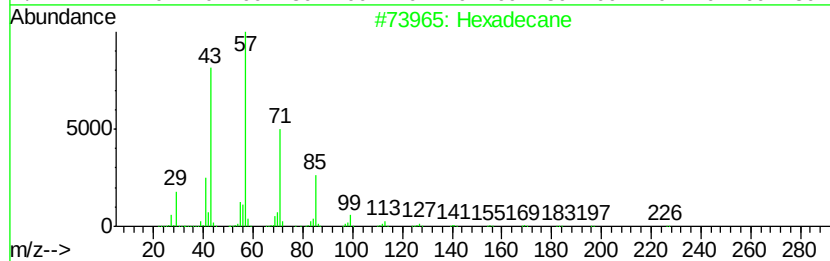
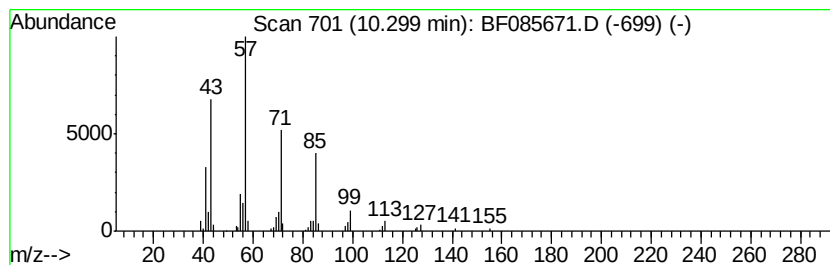
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Hexadecane Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.30	2.16 ng	123844	Acenaphthene-d10	9.86

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane		226	C16H34	000544-76-3	90
2	Pentadecane		212	C15H32	000629-62-9	90
3	Eicosane		282	C20H42	000112-95-8	83
4	Decane, 2,3,5-trimethyl-		184	C13H28	062238-11-3	83
5	Tetradecane		198	C14H30	000629-59-4	80



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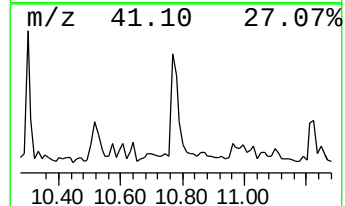
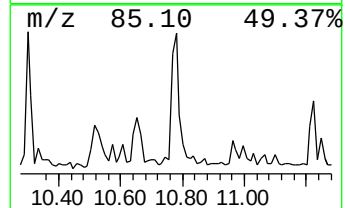
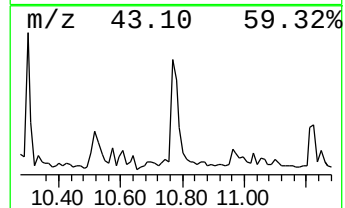
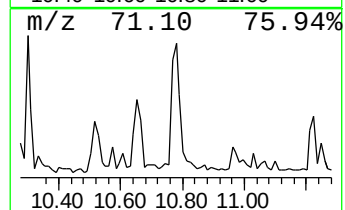
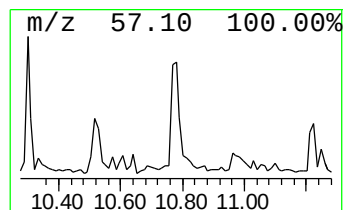
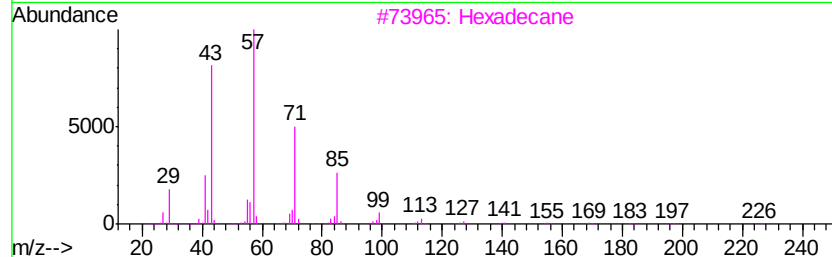
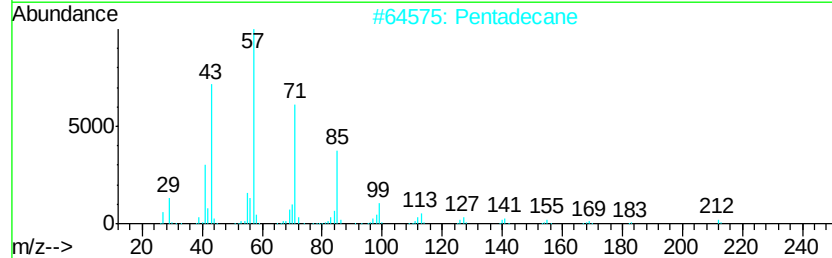
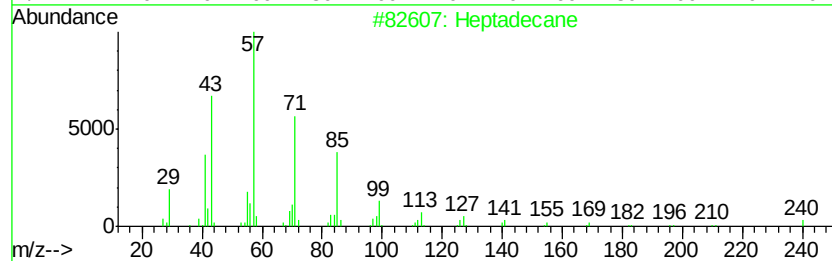
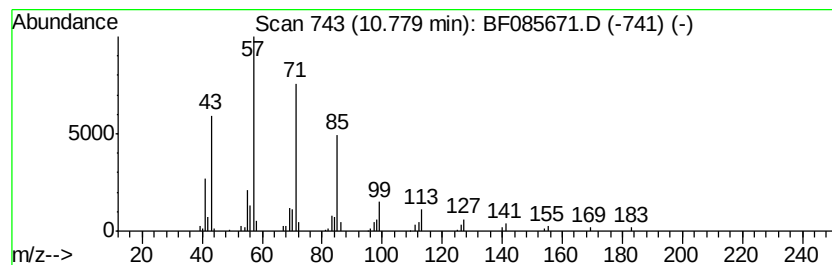
Instrument :
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 6 Heptadecane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD		R.T.
10.78	2.95 ng	188537	Phenanthrene-d10		11.35
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane	240	C17H36	000629-78-7	91
2	Pentadecane	212	C15H32	000629-62-9	91
3	Hexadecane	226	C16H34	000544-76-3	90
4	Octadecane	254	C18H38	000593-45-3	90
5	Eicosane	282	C20H42	000112-95-8	90



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ALS Vial : 9 Sample Multiplier: 1

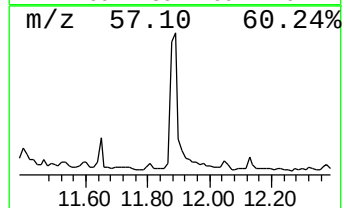
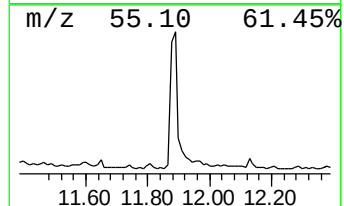
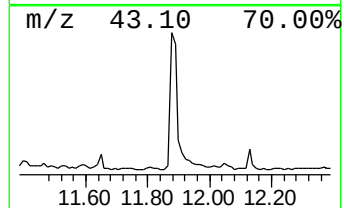
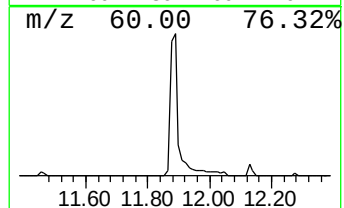
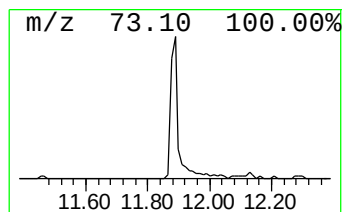
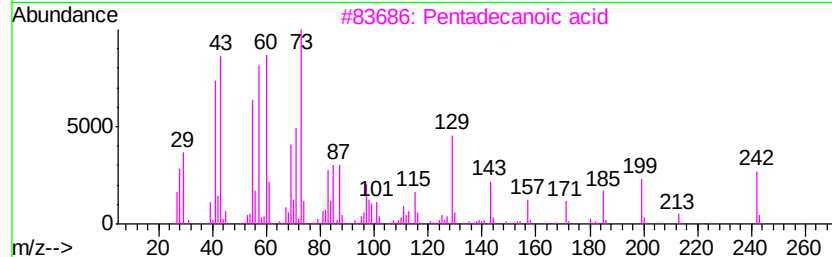
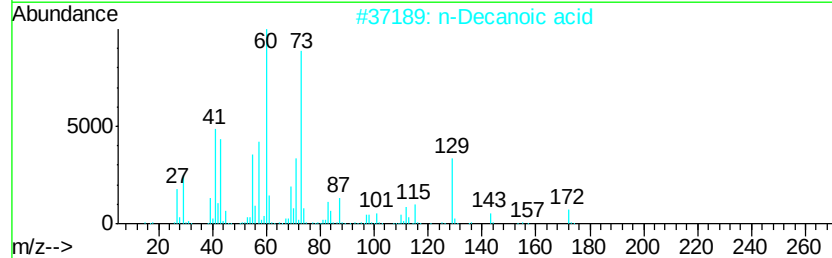
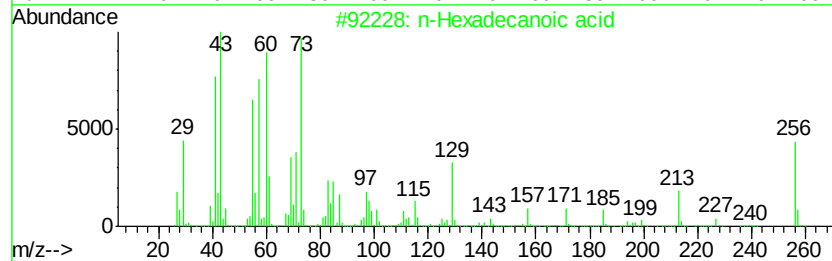
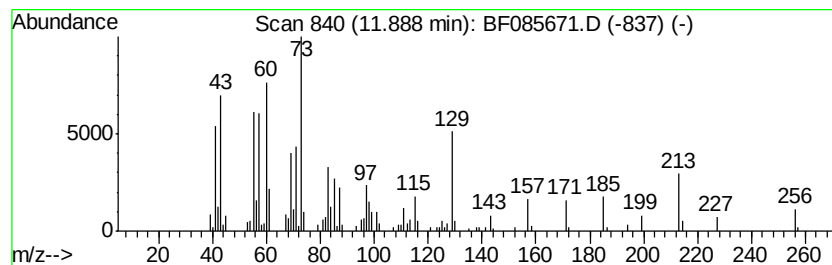
Instrument :
BNA_F
ClientSampleId :
SB-01(23-25)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031816.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 n-Hexadecanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD		R.T.	
11.89	5.15 ng	329265	Phenanthrene-d10		11.35	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid		256	C16H32O2	000057-10-3	97
2	n-Decanoic acid		172	C10H20O2	000334-48-5	76
3	Pentadecanoic acid		242	C15H30O2	001002-84-2	72
4	Tridecanoic acid		214	C13H26O2	000638-53-9	68
5	Tetradecanoic acid		228	C14H28O2	000544-63-8	52



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032216\
Data File : BF085671.D
Acq On : 22 Mar 2016 21:08
Operator : UM/SJ
Sample : H1857-02
Misc :
ALS Vial : 9 Sample Multiplier: 1

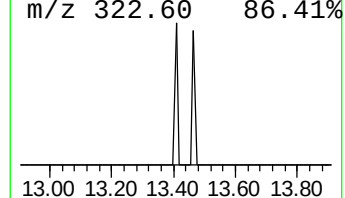
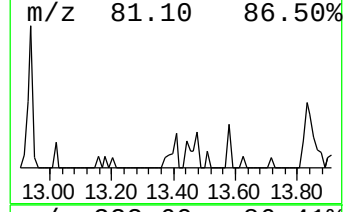
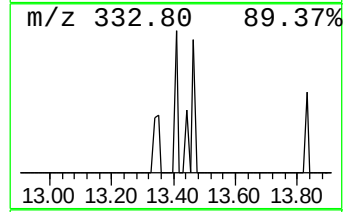
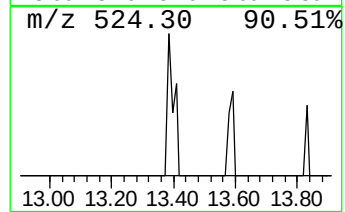
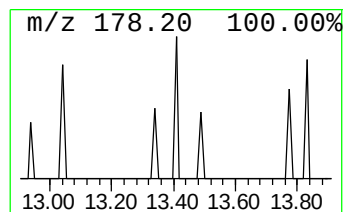
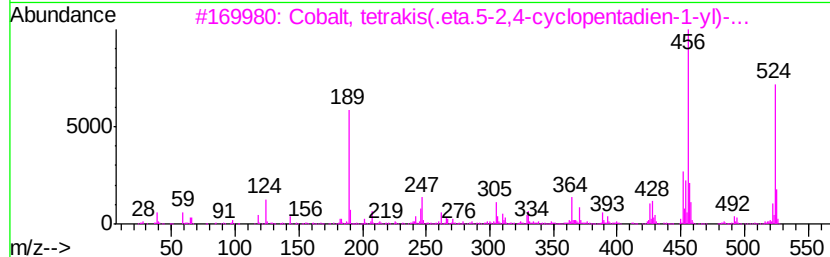
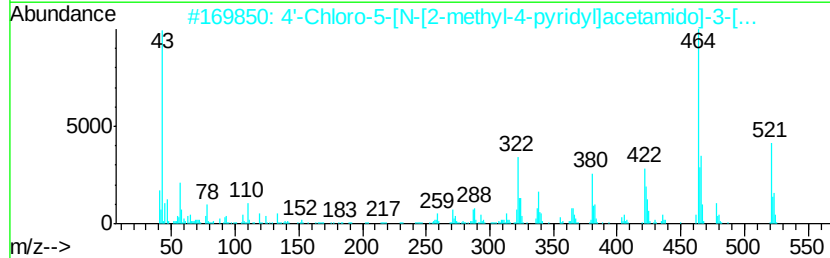
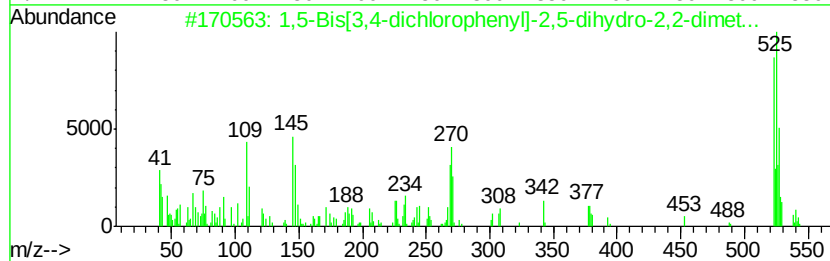
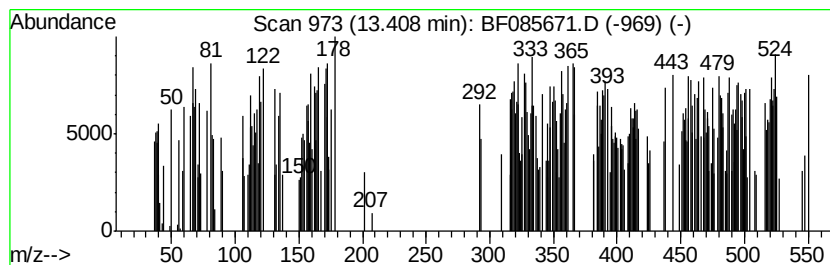
Instrument :
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ClientSampleId :
SB-01(23-25)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031816.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 1,5-Bis[3,4-dichlorophenyl]... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.		
13.41	5.13 ng	290070	Chrysene-d12	13.98		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,5-Bis[3,4-dichlorophenyl]-2,5-...	538	C27H18Cl4N4	1000212-73-4	59	
2	4'-Chloro-5-[N-[2-methyl-4-pyrid...	521	C29H32ClN3O4	1000213-37-4	10	
3	Cobalt, tetrakis(.eta.5-2,4-cycl...	524	C22H24Co4	105187-37-9	9	
4	1:12,2:3,4:5,6:7,8:9,10:11-Hexab...	522	C42H18	000190-24-9	9	
5	trans-4'-Methylthio-4-(octadecyl...	522	C34H50O2S	131316-23-9	9	



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032216\
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Acq On : 22 Mar 2016 21:08
Operator : UM/SJ
Sample : H1857-02
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-01(23-25)

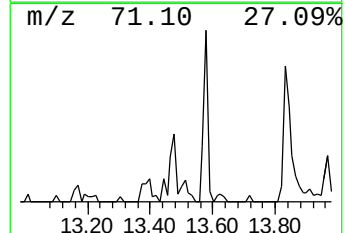
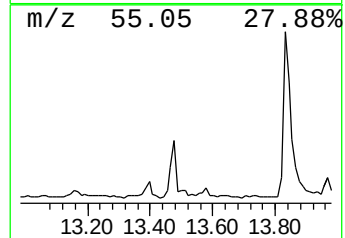
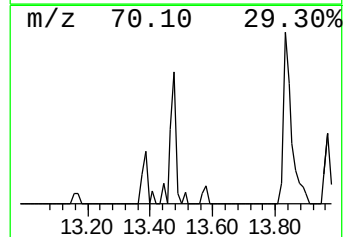
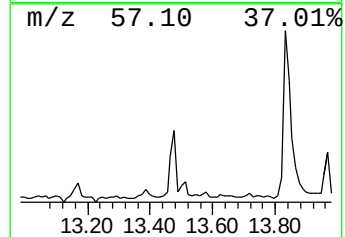
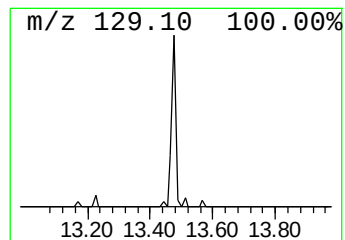
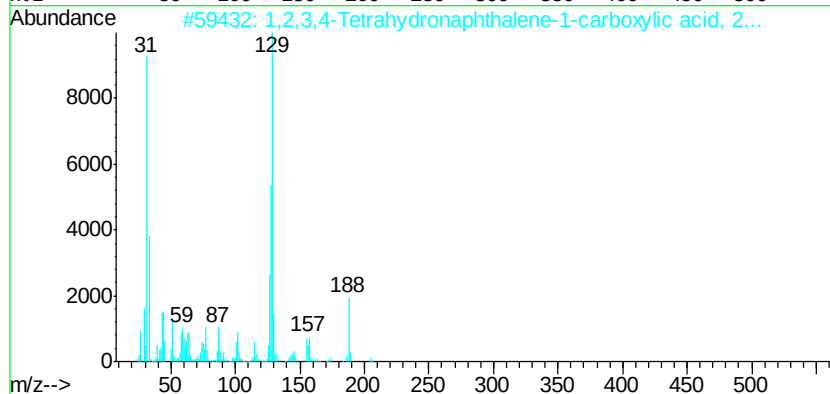
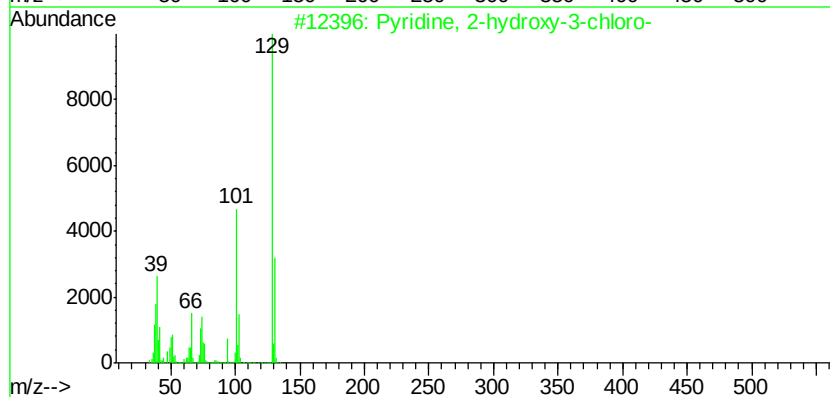
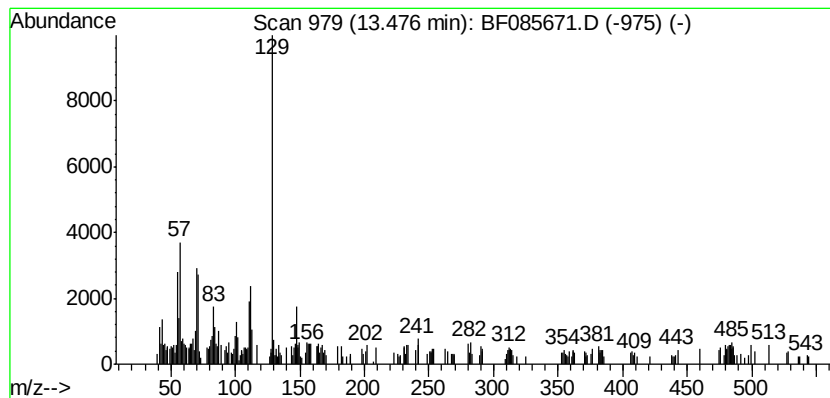
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 unknown13.48 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.48	6.81 ng	385062	Chrysene-d12	13.98

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	2	Pyridine, 2-hydroxy-3-chloro-	129	C5H4ClNO	013466-35-8	38
2		1,2,3,4-Tetrahydronaphthalene-1-...	205	C12H15NO2	068157-11-9	28



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032216\
Data File : BF085671.D
Acq On : 22 Mar 2016 21:08
Operator : UM/SJ
Sample : H1857-02
Misc :
ALS Vial : 9 Sample Multiplier: 1

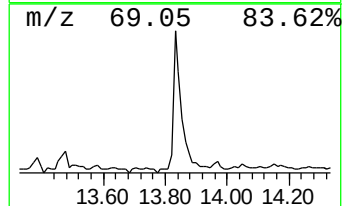
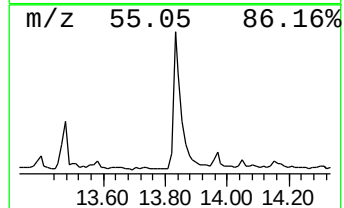
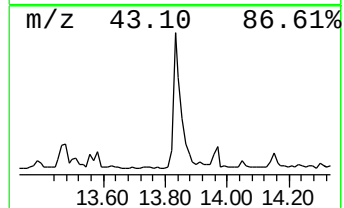
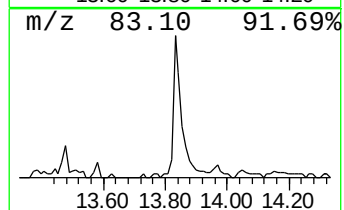
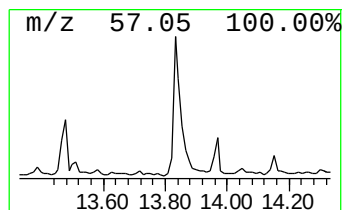
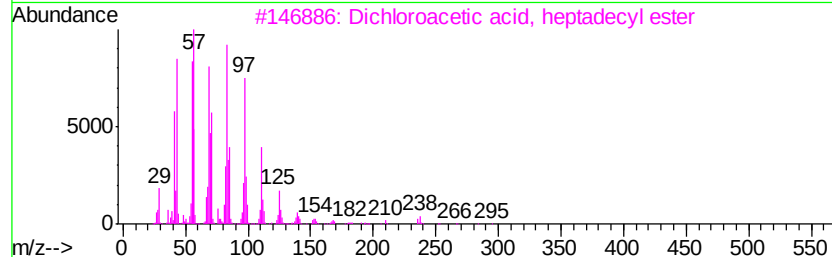
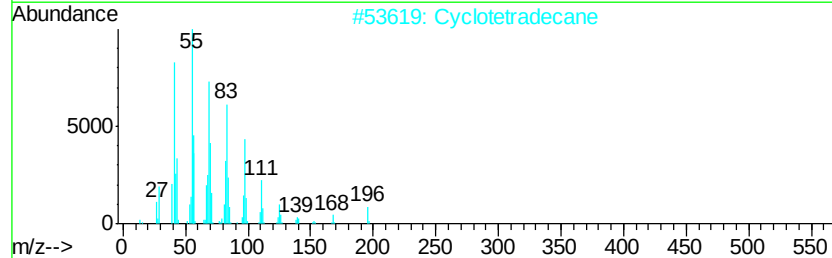
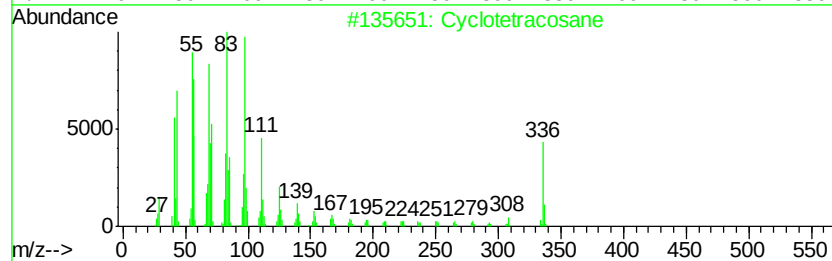
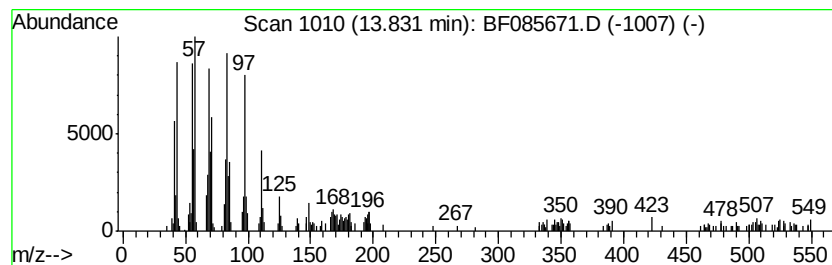
Instrument :
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ClientSampleId :
SB-01(23-25)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031816.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Cyclotetracosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD		R.T.
13.83	6.68 ng	377727	Chrysene-d12		13.98
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclotetracosane	336	C24H48	000297-03-0	95
2	Cyclotetradecane	196	C14H28	000295-17-0	92
3	Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	90
4	2- Chloropropionic acid, octadec...	360	C21H41ClO2	088104-31-8	78
5	Heptafluorobutanoic acid, heptad...	452	C21H35F7O2	1000282-97-3	78



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032216\
Data File : BF085671.D
Acq On : 22 Mar 2016 21:08
Operator : UM/SJ
Sample : H1857-02
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-01(23-25)

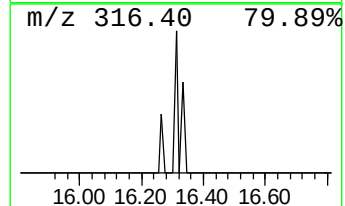
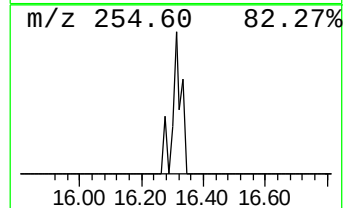
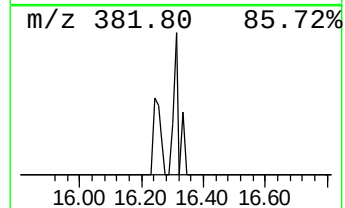
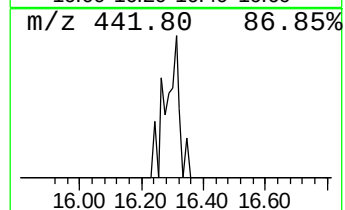
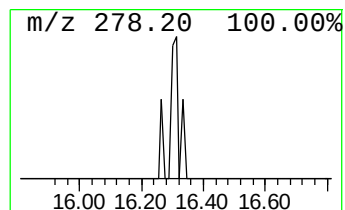
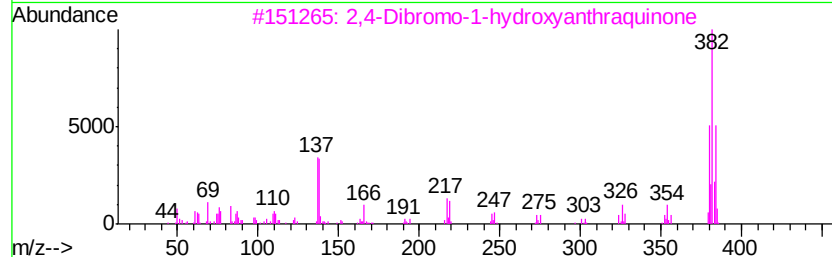
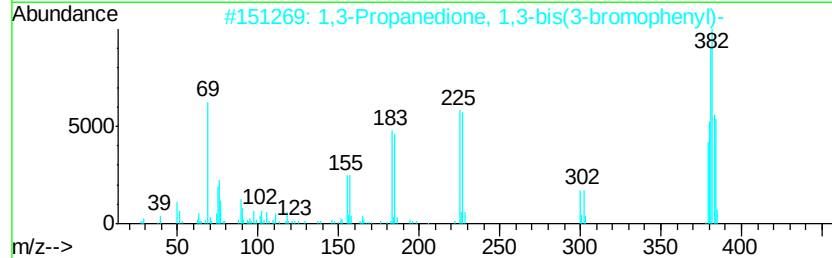
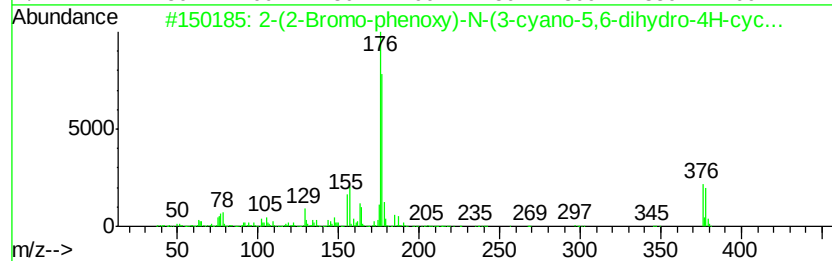
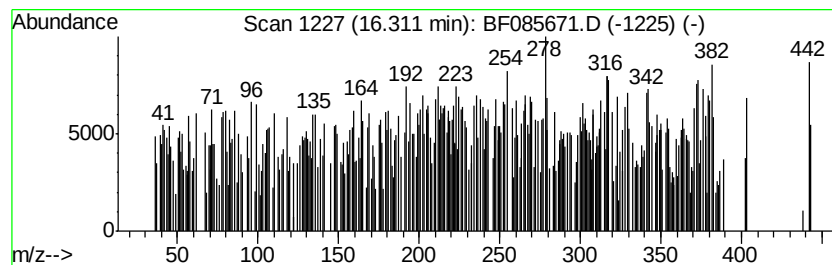
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 unknown16.31 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.31	9.17 ng	371358	Perylene-d12	15.40

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-(2-Bromo-phenoxy)-N-(3-cyano-5...	376	C16H13BrN2O2S	1000296-89-2	25
2		1,3-Propanedione, 1,3-bis(3-brom...	380	C15H10Br2O2	003387-25-5	25
3		2,4-Dibromo-1-hydroxyanthraquinone	380	C14H6Br2O3	052146-83-5	15
4		Benzene, pentachloro(trichloroet...	376	C8Cl8	029082-74-4	15
5		cis-Chlordane	406	C10H6Cl8	005103-71-9	12



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032216\
Data File : BF085671.D
Acq On : 22 Mar 2016 21:08
Operator : UM/SJ
Sample : H1857-02
Misc :
ALS Vial : 9 Sample Multiplier: 1

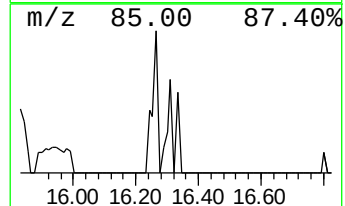
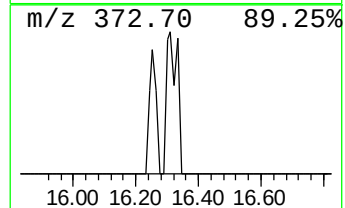
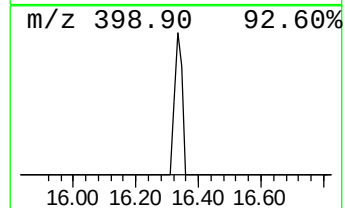
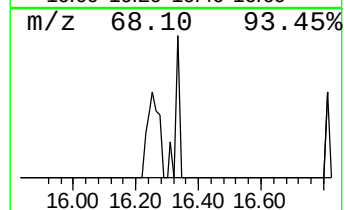
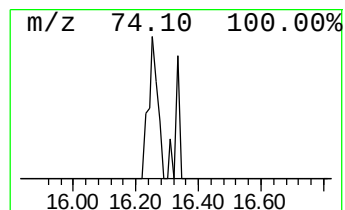
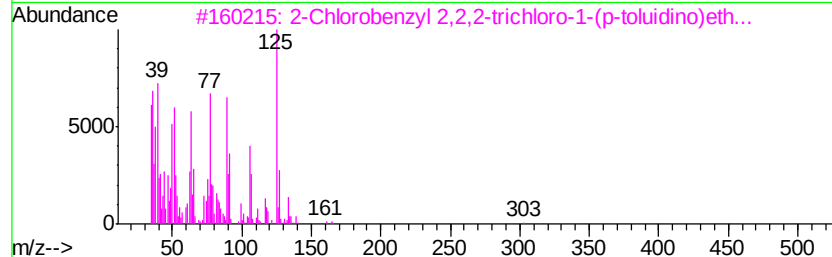
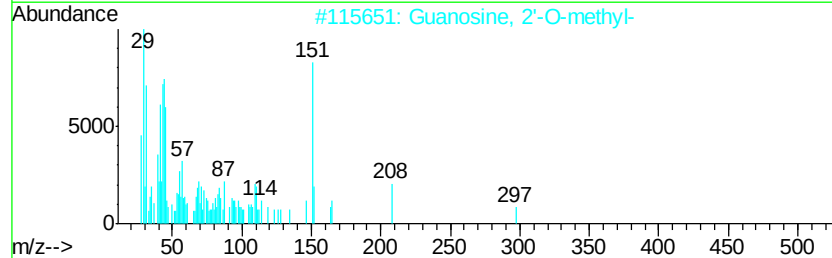
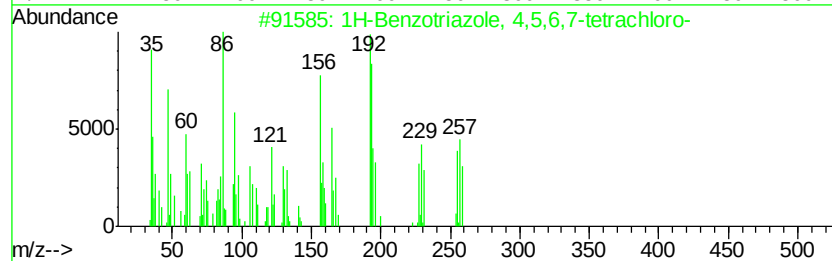
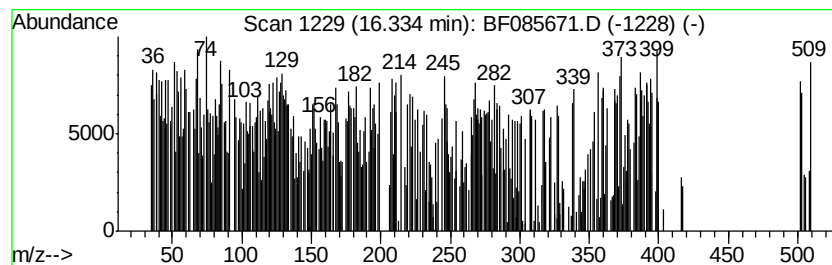
Instrument :
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ClientSampleId :
SB-01(23-25)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031816.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 1H-Benzotriazole, 4,5,6,7-t... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.		
16.33	6.38 ng	258463	Perylene-d12	15.40		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		1H-Benzotriazole, 4,5,6,7-tetrac...	255	C6HCl4N3	002338-10-5	70
2		Guanosine, 2'-O-methyl-	297	C11H15N5O5	002140-71-8	35
3		2-Chlorobenzyl 2,2,2-trichloro-1...	420	C17H16Cl4N2O2	1000222-82-6	27
4		1,2,3,4,5,8-Hexahydro-7-methoxy-...	179	C11H17NO	054893-57-1	25
5		Ethane, 1-[(2-chloroethyl)thio]-...	184	C6H13ClS2	092569-22-7	22



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032216\
 Data File : BF085671.D
 Acq On : 22 Mar 2016 21:08
 Operator : UM/SJ
 Sample : H1857-02
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031816.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
2-Pentanone, 4-hy...	5.02	31.4	ng	1164570	1	6.84	741167	20.0
2-Pentanone, 4-me...	5.84	3.0	ng	109668	1	6.84	741167	20.0
unknown6.60	6.60	101.1	ng	3744790	1	6.84	741167	20.0
Hexadecane	10.30	2.2	ng	123844	3	9.86	1148730	20.0
Heptadecane	10.78	3.0	ng	188537	4	11.35	1278140	20.0
n-Hexadecanoic acid	11.89	5.2	ng	329265	4	11.35	1278140	20.0
1,5-Bis[3,4-dichl...	13.41	5.1	ng	290070	5	13.98	1131310	20.0
unknown13.48	13.48	6.8	ng	385062	5	13.98	1131310	20.0
Cyclotetracosane	13.83	6.7	ng	377727	5	13.98	1131310	20.0
unknown16.27	16.27	24.8	ng	1005000	6	15.40	810064	20.0
unknown16.31	16.31	9.2	ng	371358	6	15.40	810064	20.0
1H-Benzotriazole,...	16.33	6.4	ng	258463	6	15.40	810064	20.0