

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050715\
 Data File : BF079006.D
 Acq On : 7 May 2015 20:07
 Operator : TP/IZ
 Sample : G2084-01 5X
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 1

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-BF043015.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	1.484	23	26	28	rVB	860209	1178650	18.80%	6.699%
2	1.621	36	38	48	rVB2	125043	323403	5.16%	1.838%
3	1.758	48	50	52	rVV	96439	152697	2.44%	0.868%
4	1.804	52	54	60	rVV	69176	150291	2.40%	0.854%
5	1.884	60	61	98	rVB	2732631	6269098	100.00%	35.632%
6	5.644	387	390	392	rVB	197963	260747	4.16%	1.482%
7	6.902	497	500	502	rBV	228272	279533	4.46%	1.589%
8	7.050	511	513	516	rVB	320745	300173	4.79%	1.706%
9	7.336	536	538	540	rBV	278193	360591	5.75%	2.050%
10	7.530	553	555	557	rVB	207289	207216	3.31%	1.178%
11	8.033	596	599	601	rVB	146572	155127	2.47%	0.882%
12	8.913	674	676	679	rBV	410367	506901	8.09%	2.881%
13	10.262	792	794	796	rVB	316472	268706	4.29%	1.527%
14	11.096	864	867	870	rVV	647216	612455	9.77%	3.481%
15	11.668	913	917	920	rBV5	31350	79027	1.26%	0.449%
16	11.759	923	925	929	rVV	86514	94793	1.51%	0.539%
17	11.942	938	941	943	rBV	134926	215794	3.44%	1.227%
18	12.079	950	953	955	rBV	244718	286836	4.58%	1.630%
19	12.182	960	962	965	rVB	419607	402645	6.42%	2.289%
20	12.285	968	971	973	rBV	358887	517101	8.25%	2.939%
21	12.514	988	991	992	rBV	81101	103153	1.65%	0.586%
22	12.548	992	994	995	rVB	66721	64564	1.03%	0.367%
23	12.868	1019	1022	1024	rVB	624653	903591	14.41%	5.136%
24	12.948	1026	1029	1031	rBV	640613	783993	12.51%	4.456%
25	13.005	1032	1034	1035	rVV	90784	88973	1.42%	0.506%
26	13.302	1058	1060	1062	rBV	310698	373800	5.96%	2.125%
27	13.417	1068	1070	1072	rVB	107929	150712	2.40%	0.857%
28	14.194	1136	1138	1139	rBV	205141	235821	3.76%	1.340%
29	14.457	1159	1161	1163	rBV	117758	166080	2.65%	0.944%
30	14.491	1163	1164	1166	rVB	119910	111822	1.78%	0.636%
31	14.948	1201	1204	1205	rBV	211200	311620	4.97%	1.771%
32	16.274	1317	1320	1322	rBV	680800	839663	13.39%	4.772%
33	18.137	1481	1483	1486	rBV	543763	838380	13.37%	4.765%

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Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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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-BF043015.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

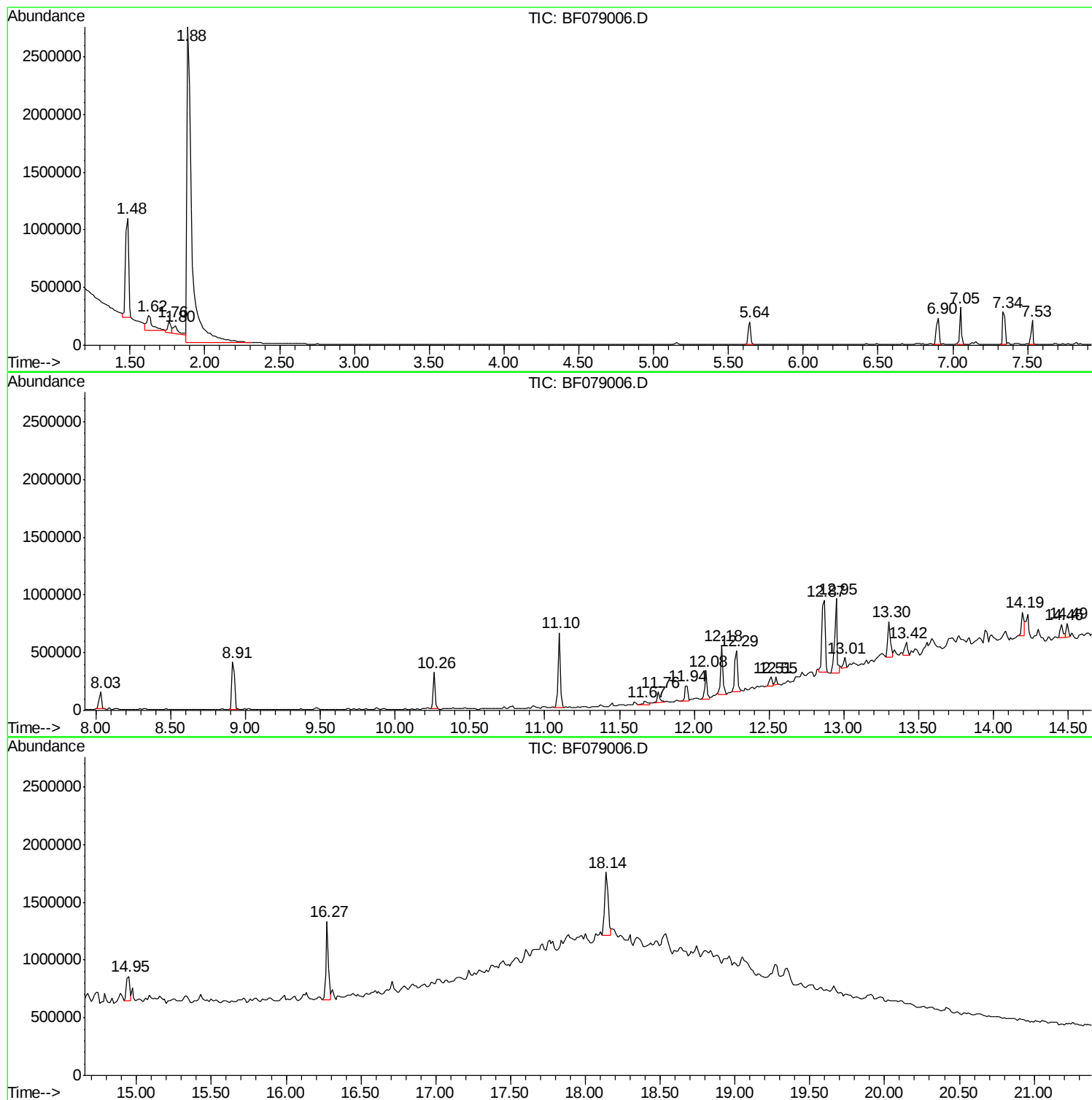
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF043015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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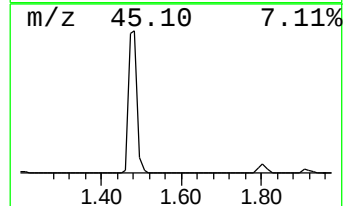
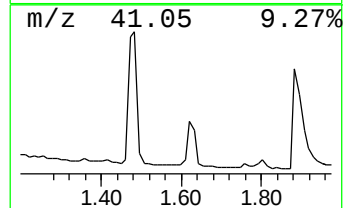
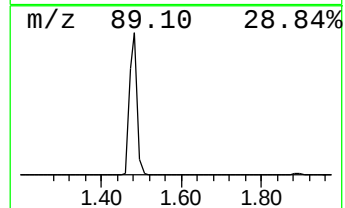
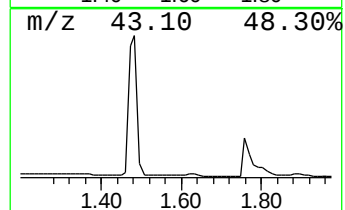
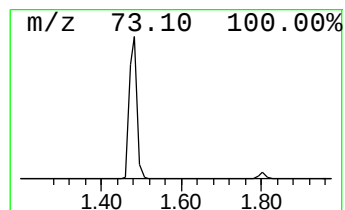
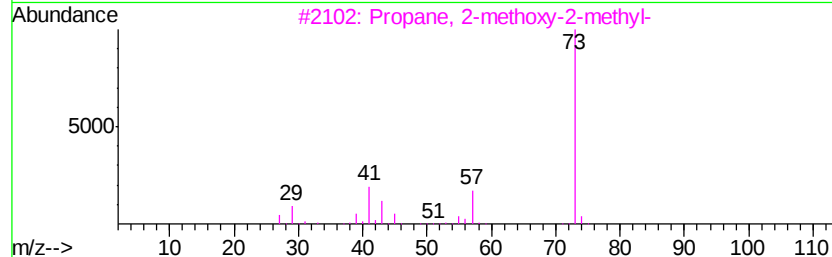
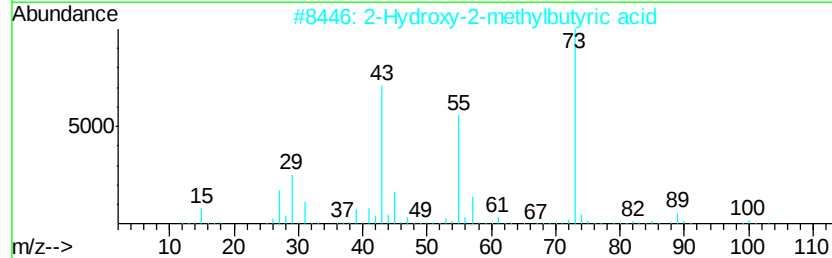
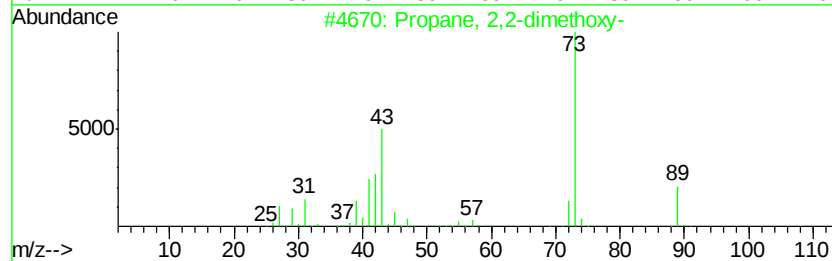
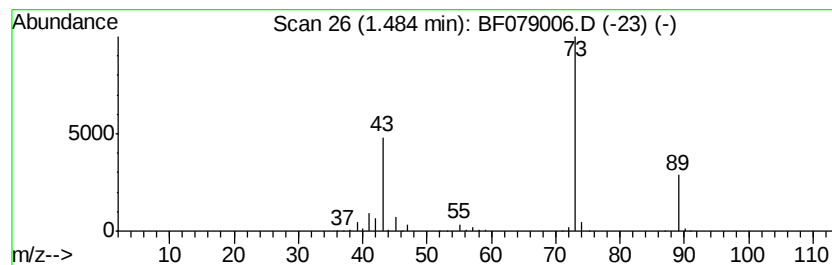
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Propane, 2,2-dimethoxy- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.48	65.37 ng	1178650	1,4-Dichlorobenzene-d4	7.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propane, 2,2-dimethoxy-	104	C5H12O2	000077-76-9	56
2		2-Hydroxy-2-methylbutyric acid	118	C5H10O3	003739-30-8	9
3		Propane, 2-methoxy-2-methyl-	88	C5H12O	001634-04-4	9
4		Methane, isothiocyanato-	73	C2H3NS	000556-61-6	7
5		Borane, dimethoxy-	74	C2H7BO2	004542-61-4	7



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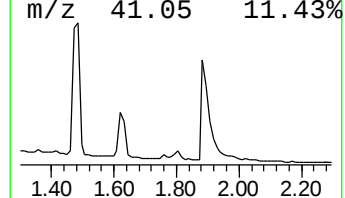
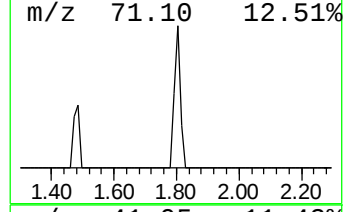
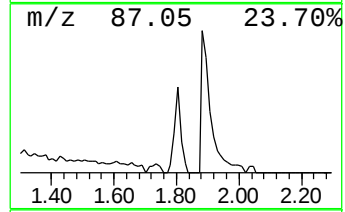
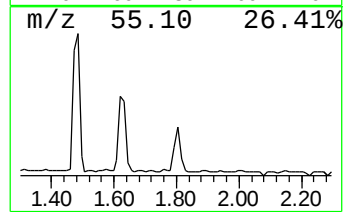
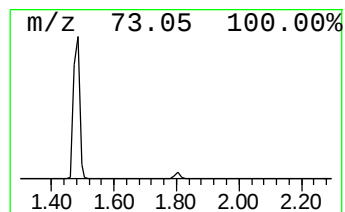
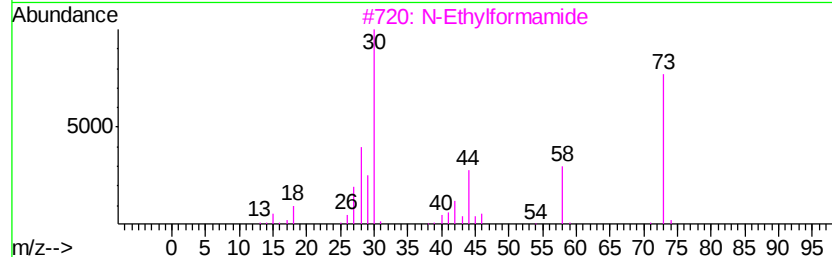
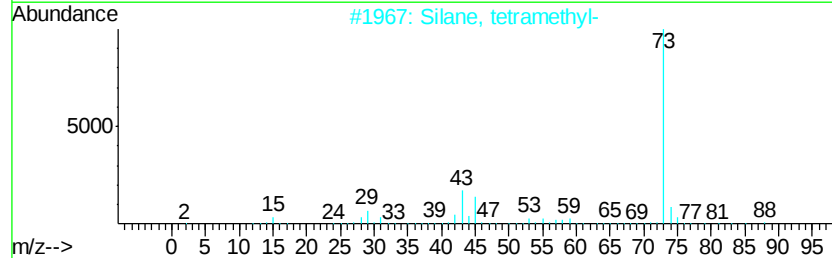
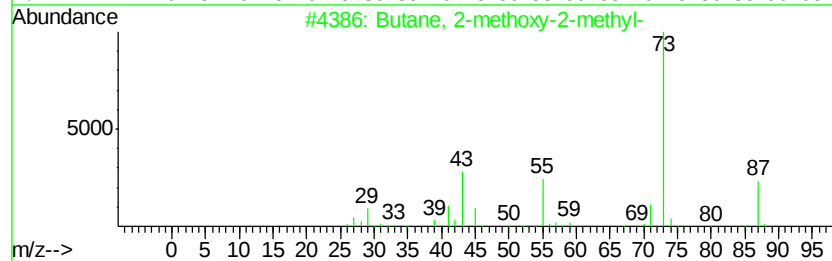
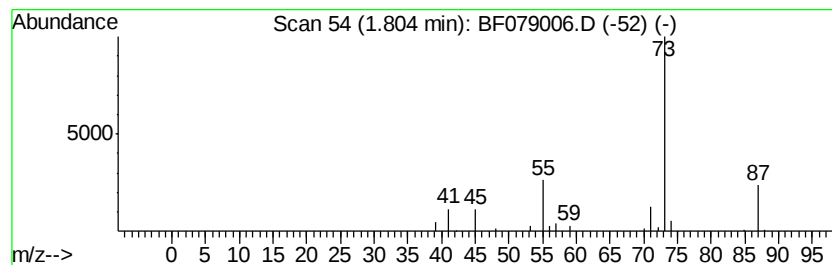
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Butane, 2-methoxy-2-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.80	8.34 ng	150291	1,4-Dichlorobenzene-d4	7.35

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	64
2		Silane, tetramethyl-	88	C4H12Si	000075-76-3	32
3		N-Ethylformamide	73	C3H7NO	000627-45-2	9
4		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	9
5		Silane, trimethyl[2-methylene-1-...	194	C12H22Si	097778-15-9	9



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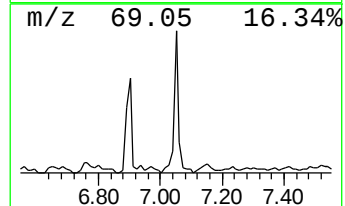
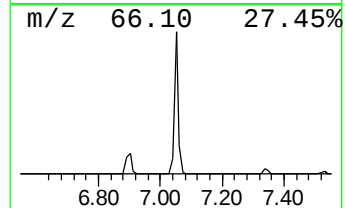
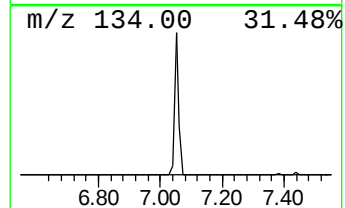
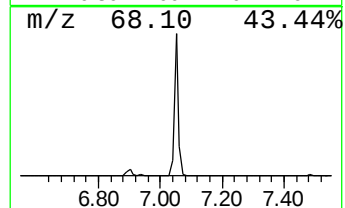
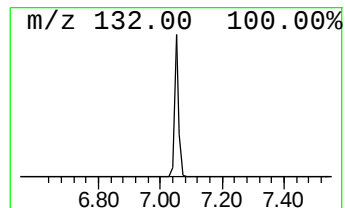
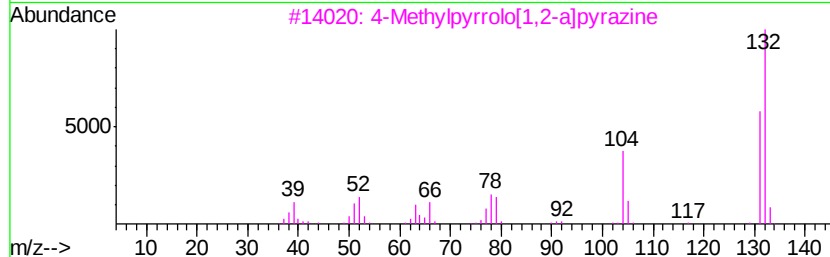
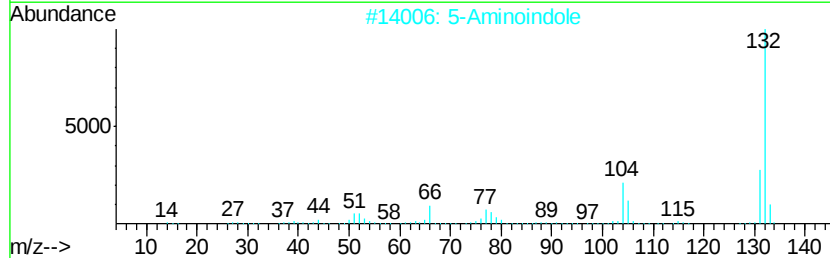
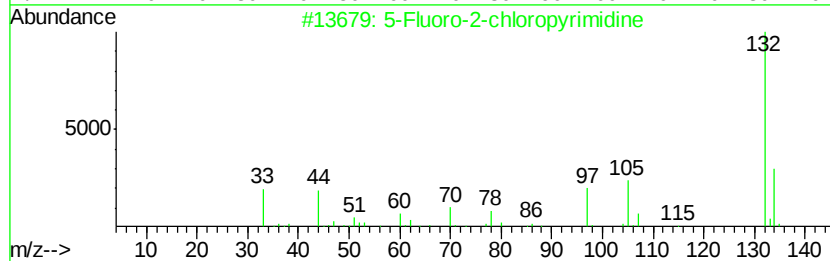
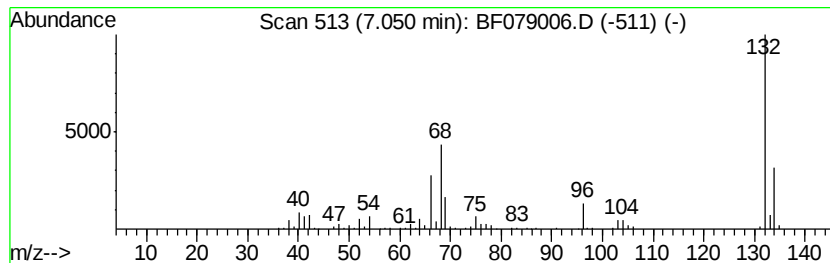
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 unknown7.05 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.05	16.65 ng	300173	1,4-Dichlorobenzene-d4	7.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
2		5-Aminoindole	132	C8H8N2	005192-03-0	10
3		4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9
4		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9
5		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9



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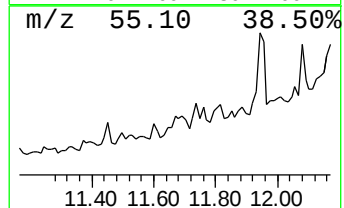
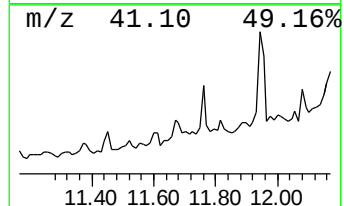
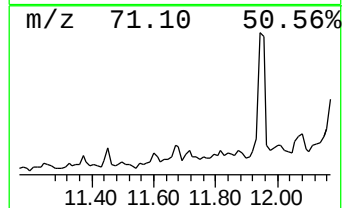
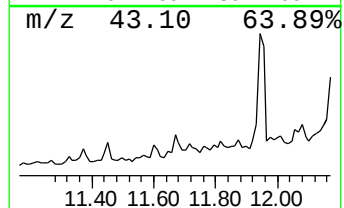
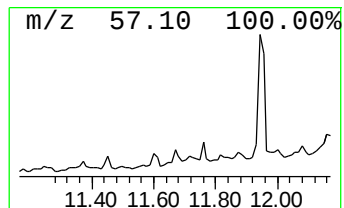
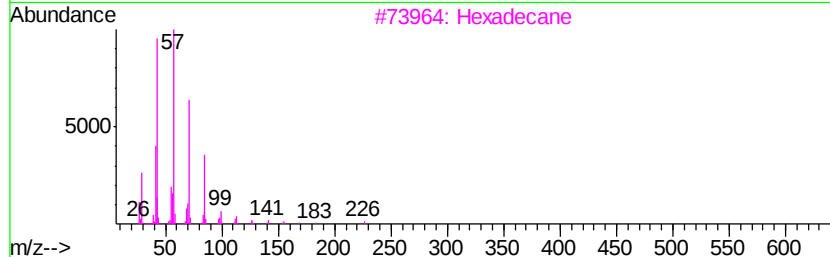
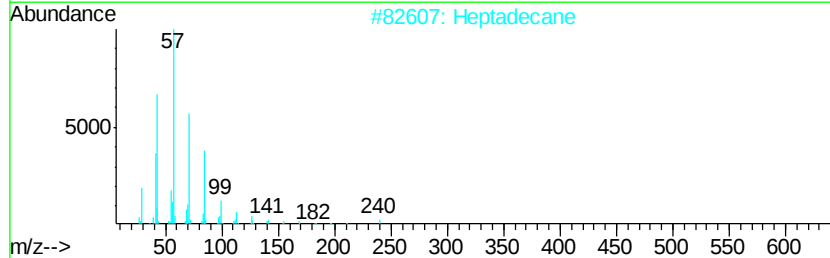
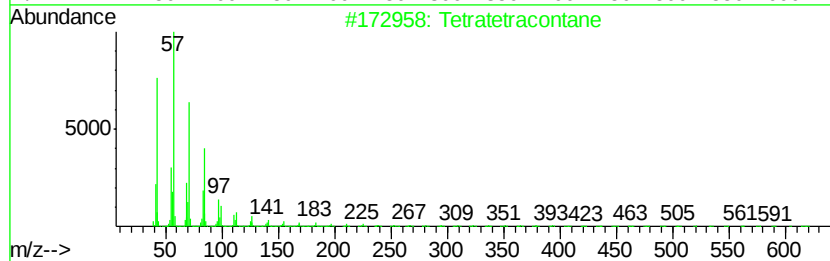
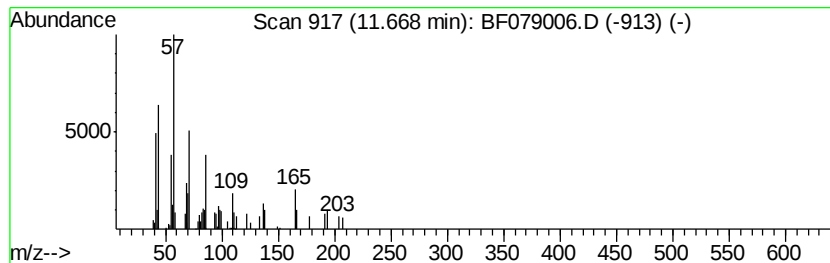
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 unknown11.67 Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.67	2.58 ng	79027	Acenaphthene-d10	11.10

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetratetracontane	619	C44H90	007098-22-8	38
2			Heptadecane	240	C17H36	000629-78-7	35
3			Hexadecane	226	C16H34	000544-76-3	35
4			Undecane, 5-ethyl-	184	C13H28	017453-94-0	35
5			Dodecane	170	C12H26	000112-40-3	35



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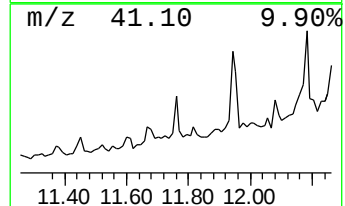
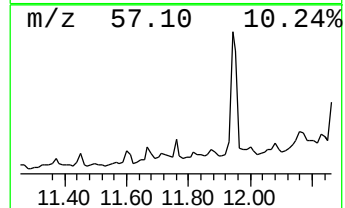
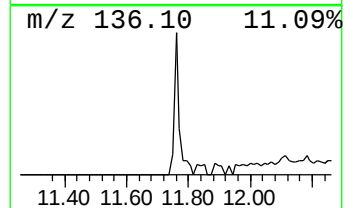
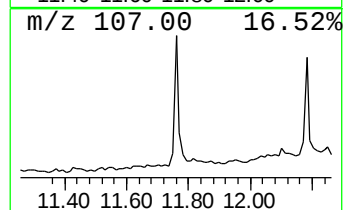
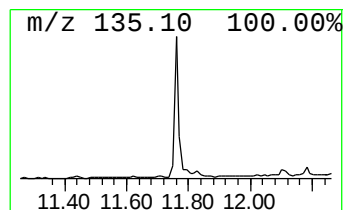
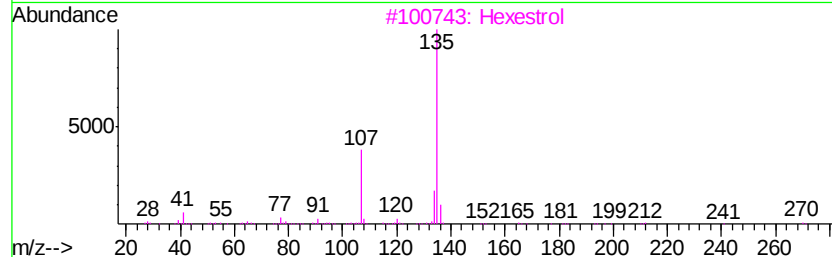
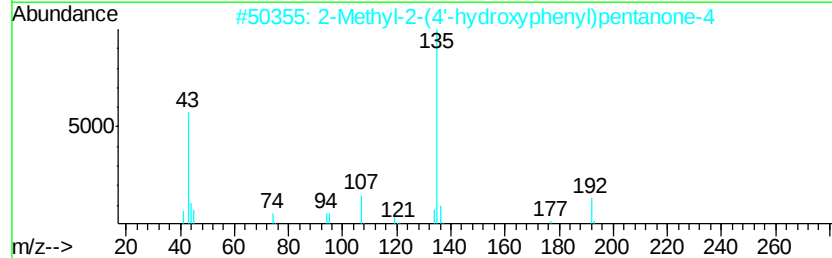
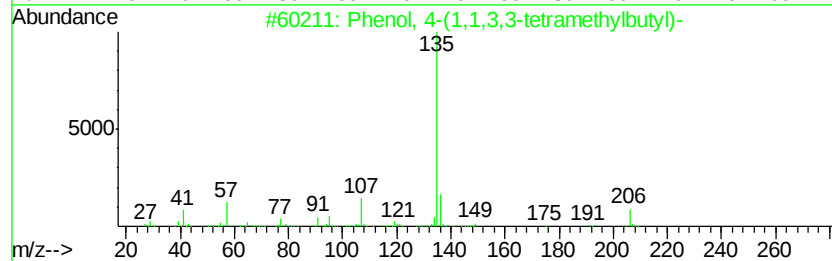
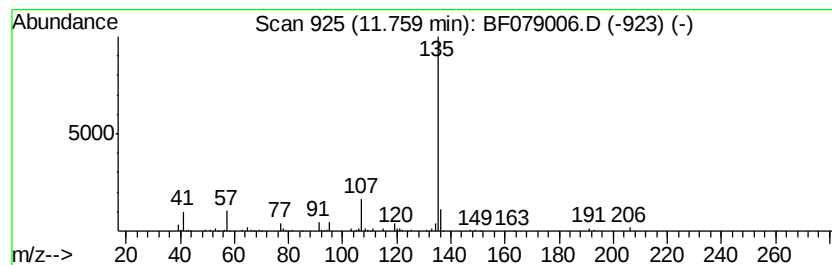
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Phenol, 4-(1,1,3,3-tetramet... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.76	3.10 ng	94793	Acenaphthene-d10	11.10	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 4-(1,1,3,3-tetramethylbu...	206	C14H22O	000140-66-9	91
2	2-Methyl-2-(4'-hydroxyphenyl)pen...	192	C12H16O2	070205-18-4	64
3	Hexestrol	270	C18H22O2	005635-50-7	59
4	Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	56
5	1-Adamantan-1-yl-6-chloro-7-fluo...	375	C20H19ClFN03	1000210-85-1	53



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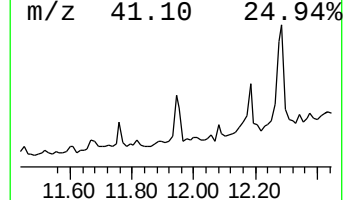
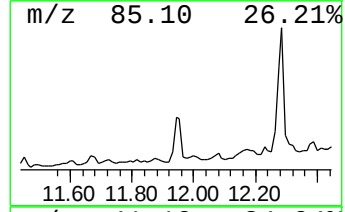
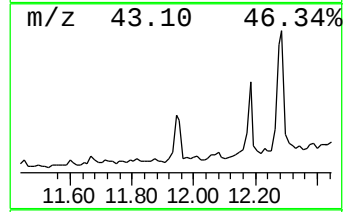
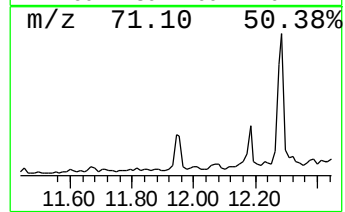
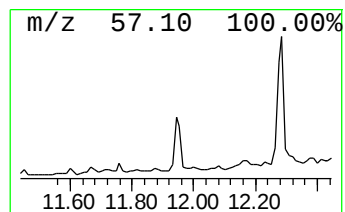
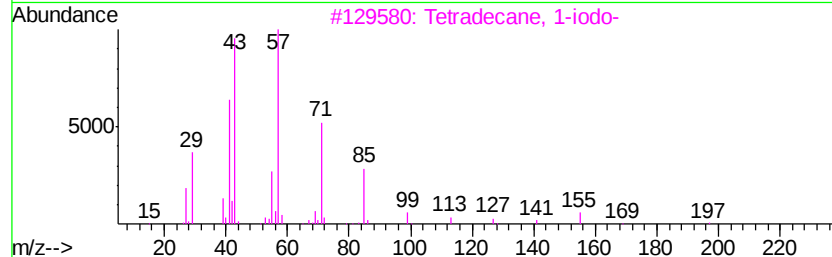
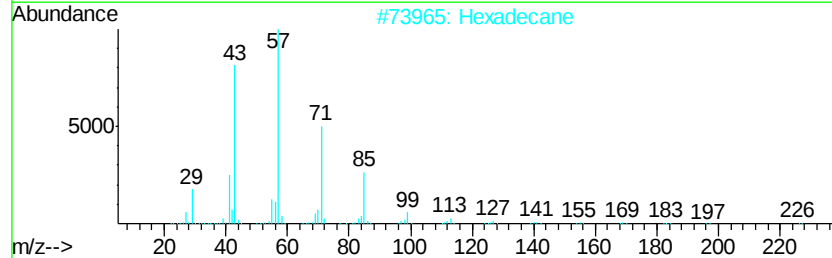
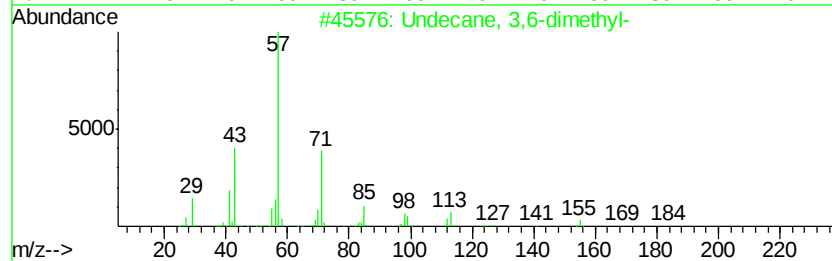
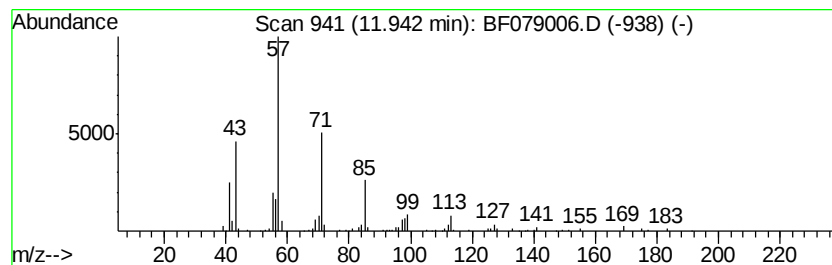
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Undecane, 3,6-dimethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.94	7.05 ng	215794	Acenaphthene-d10	11.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Undecane, 3,6-dimethyl-	184	C13H28	017301-28-9	72
2		Hexadecane	226	C16H34	000544-76-3	72
3		Tetradecane, 1-iodo-	324	C14H29I	019218-94-1	59
4		Octadecane	254	C18H38	000593-45-3	58
5		Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	58



Data Path : Z:\HPCHEM1\BNA F\DATA\BF050715\
Data File : BF079006.D
Acq On : 7 May 2015 20:07
Operator : TP/IZ
Sample : G2084-01 5X
Misc :
ALS Vial : 17 Sample Multiplier: 1

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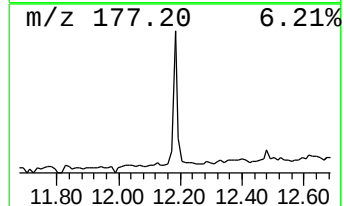
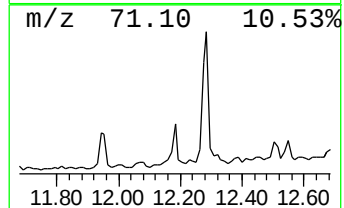
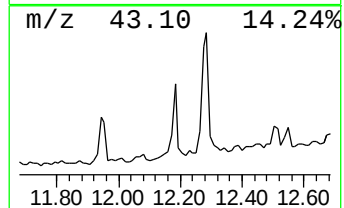
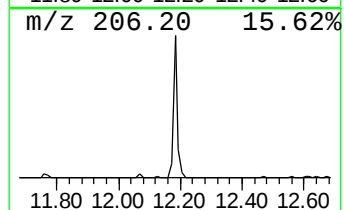
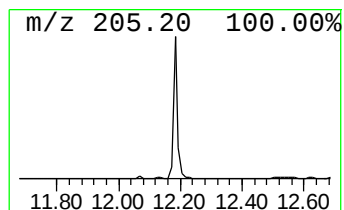
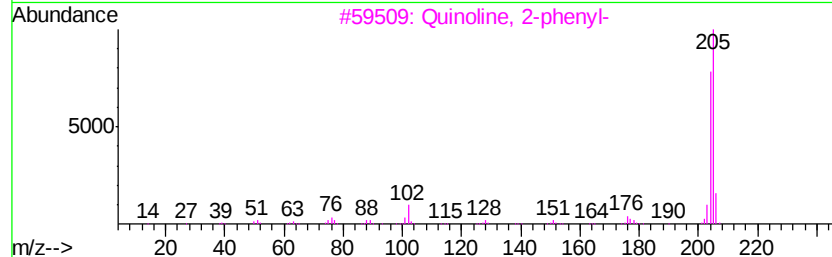
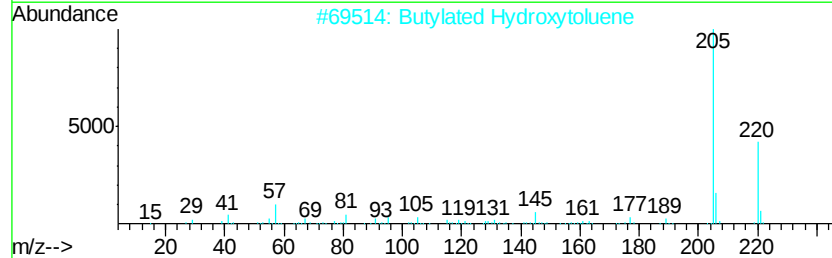
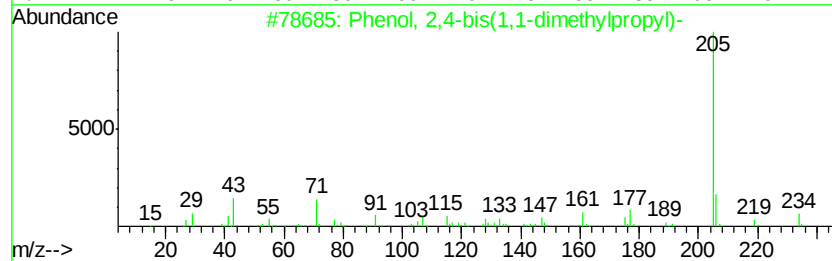
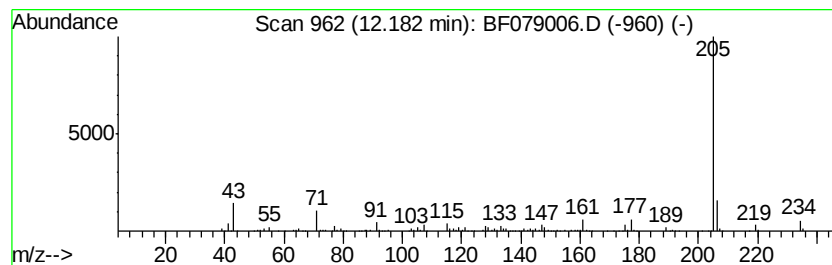
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Phenol, 2,4-bis(1,1-dimethy... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.18	10.27 ng	402645	Phenanthrene-d10	12.95

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenol, 2,4-bis(1,1-dimethylprop...	234	C16H26O	000120-95-6	92
2		Butylated Hydroxytoluene	220	C15H24O	000128-37-0	59
3		Quinoline, 2-phenyl-	205	C15H11N	000612-96-4	59
4		Phenol, 2,6-bis(1,1-dimethylethyl...	277	C17H27NO2	001918-11-2	59
5		6-Phenylisoquinoline	205	C15H11N	070125-61-0	59



Data Path : Z:\HPCHEM1\BNA F\DATA\BF050715\
Data File : BF079006.D
Acq On : 7 May 2015 20:07
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Sample : G2084-01 5X
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ALS Vial : 17 Sample Multiplier: 1

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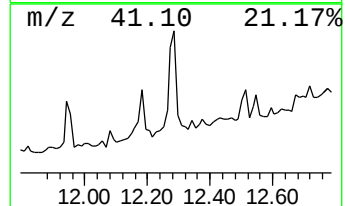
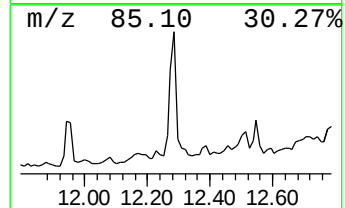
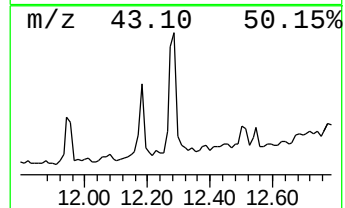
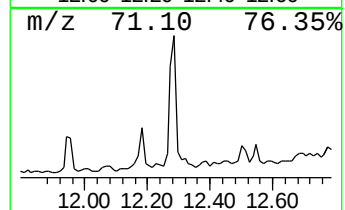
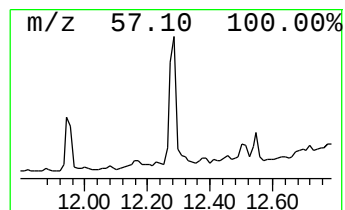
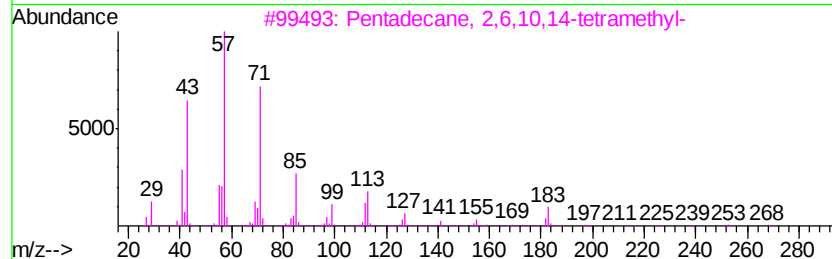
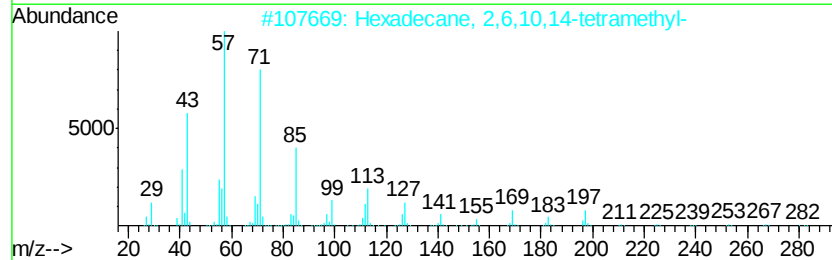
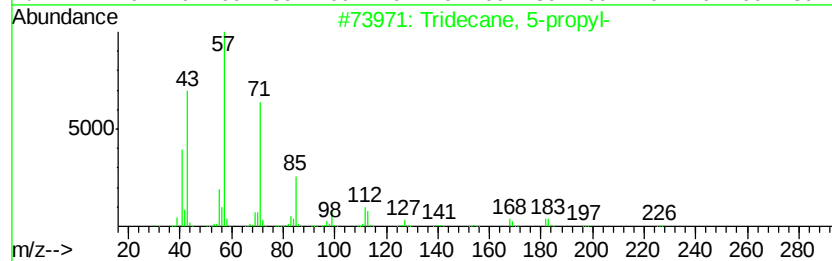
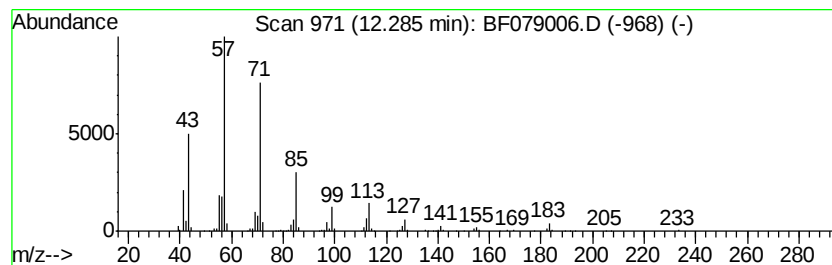
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Tridecane, 5-propyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.29	13.19 ng	517101	Phenanthrene-d10	12.95

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tridecane, 5-propyl-	226	C16H34	055045-11-9	90
2		Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	90
3		Pentadecane, 2,6,10,14-tetramethyl-	268	C19H40	001921-70-6	90
4		Dodecane, 2,6,10-trimethyl-	212	C15H32	003891-98-3	78
5		Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	72



Data Path : Z:\HPCHEM1\BNA F\DATA\BF050715\
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Sample : G2084-01 5X
Misc :
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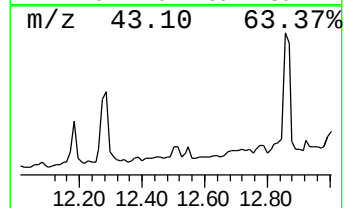
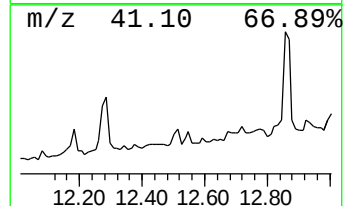
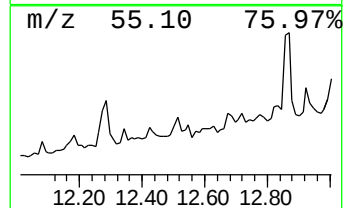
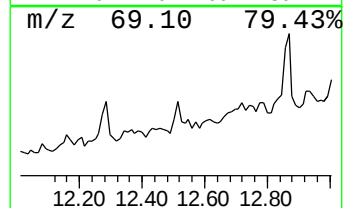
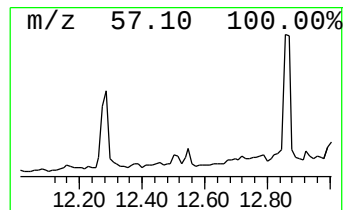
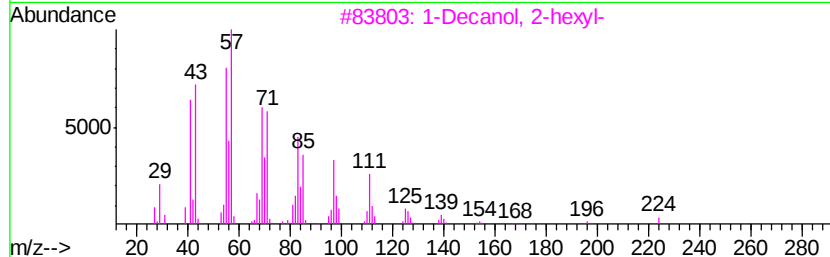
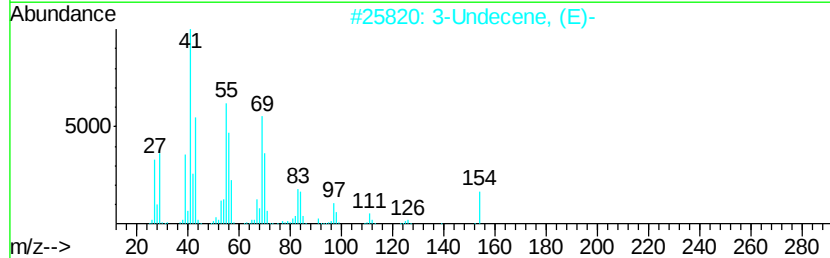
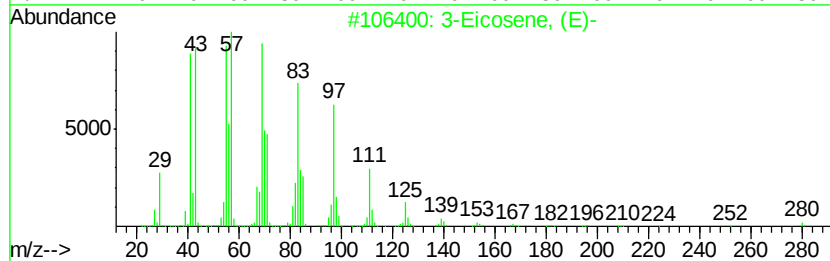
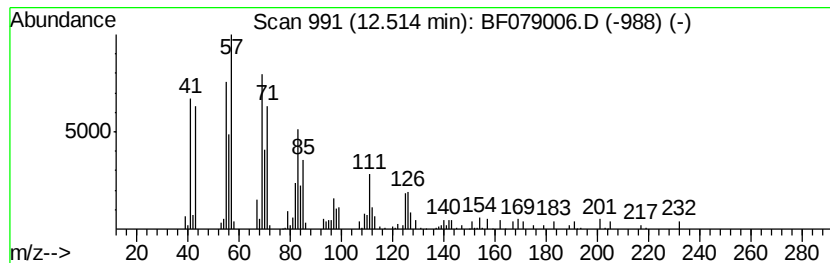
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF043015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 3-Eicosene, (E)- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.51	2.63 ng	103153	Phenanthrene-d10	12.95	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Eicosene, (E)-	280	C20H40	074685-33-9	74
2	3-Undecene, (E)-	154	C11H22	001002-68-2	64
3	1-Decanol, 2-hexyl-	242	C16H34O	002425-77-6	62
4	3-Nonene, (E)-	126	C9H18	020063-92-7	60
5	3-Undecene, (Z)-	154	C11H22	000821-97-6	60



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050715\
Data File : BF079006.D
Acq On : 7 May 2015 20:07
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Misc :
ALS Vial : 17 Sample Multiplier: 1

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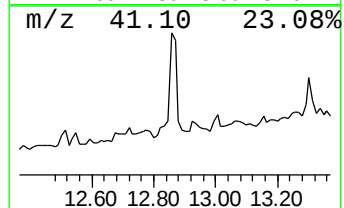
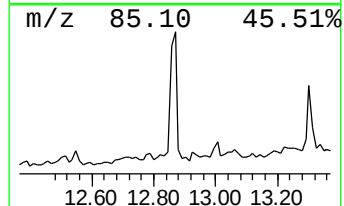
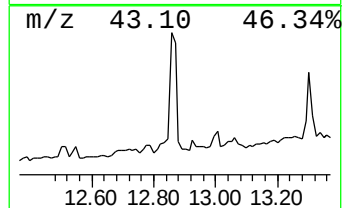
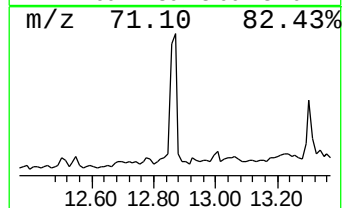
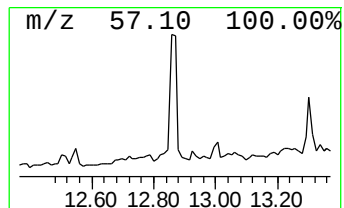
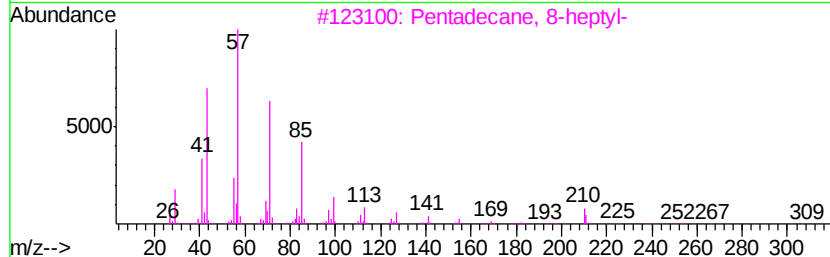
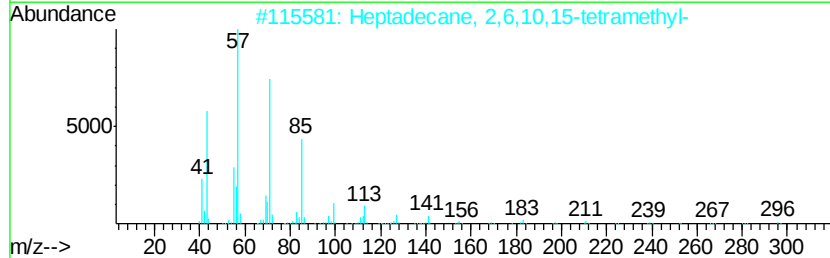
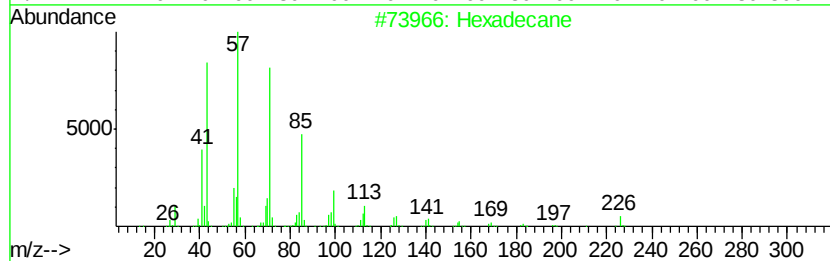
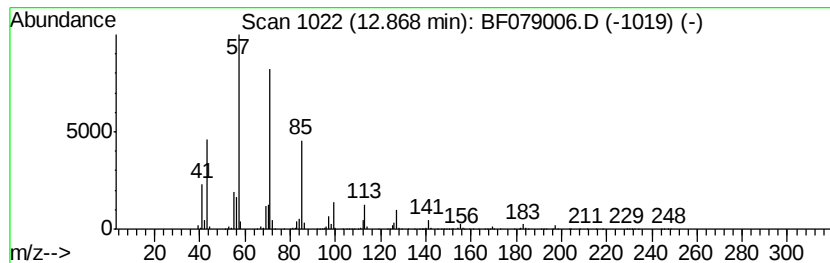
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Hexadecane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.87	23.05 ng	903591	Phenanthrene-d10	12.95

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecane	226	C16H34	000544-76-3	91
2		Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	91
3		Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	90
4		Heptadecane	240	C17H36	000629-78-7	90
5		Eicosane	282	C20H42	000112-95-8	86



Data Path : Z:\HPCHEM1\BNA F\DATA\BF050715\
Data File : BF079006.D
Acq On : 7 May 2015 20:07
Operator : TP/IZ
Sample : G2084-01 5X
Misc :
ALS Vial : 17 Sample Multiplier: 1

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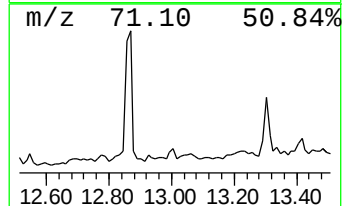
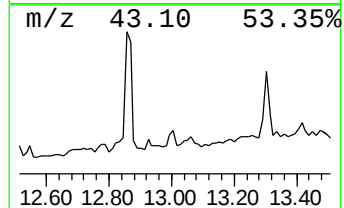
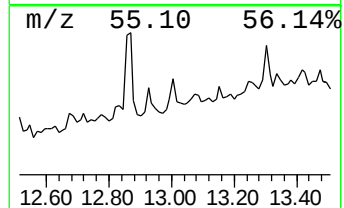
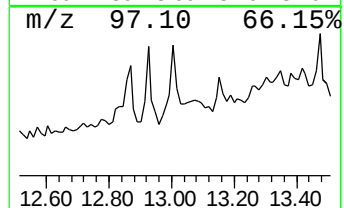
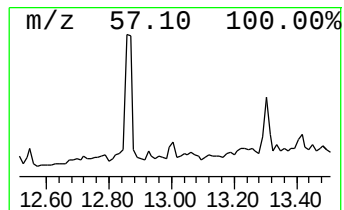
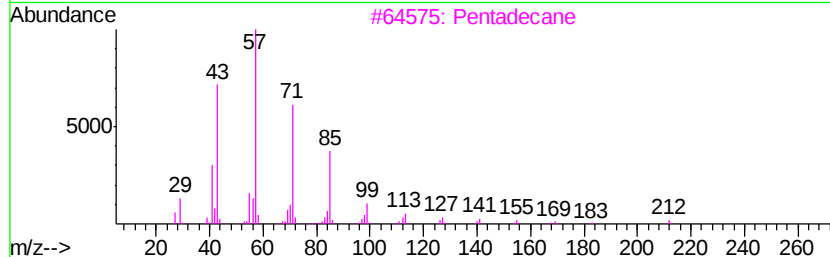
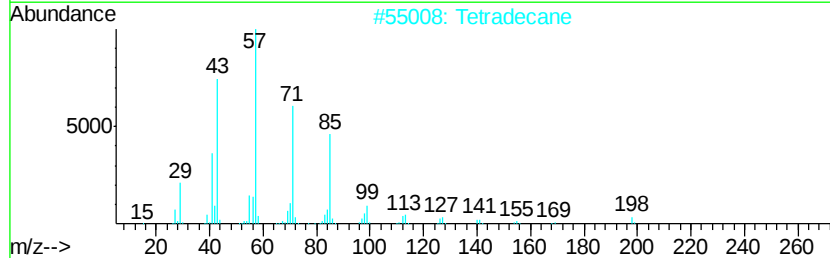
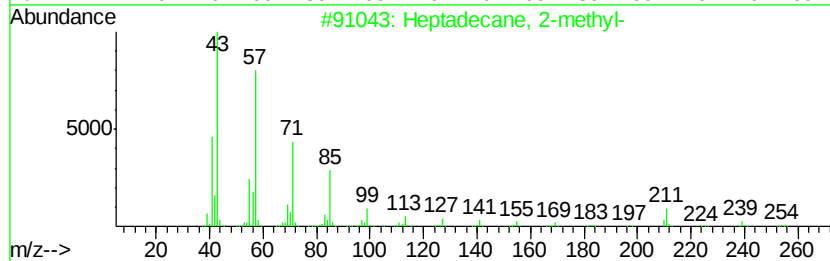
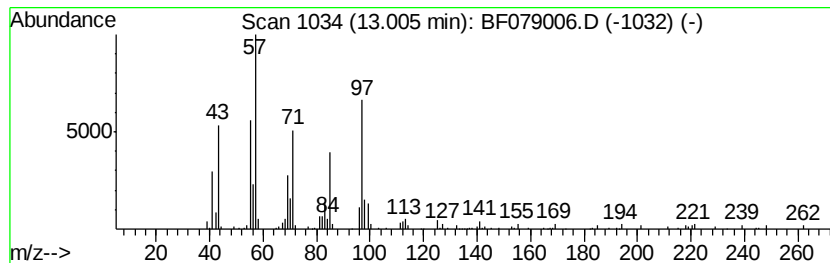
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Heptadecane, 2-methyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.01	2.27 ng	88973	Phenanthrene-d10	12.95

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane, 2-methyl-	254	C18H38	001560-89-0	53
2		Tetradecane	198	C14H30	000629-59-4	52
3		Pentadecane	212	C15H32	000629-62-9	52
4		Tetracontane, 3,5,24-trimethyl-	605	C43H88	055162-61-3	50
5		Undecane, 3,9-dimethyl-	184	C13H28	017301-31-4	50



Data Path : Z:\HPCHEM1\BNA F\DATA\BF050715\
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Acq On : 7 May 2015 20:07
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Misc :
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ClientSampleId :
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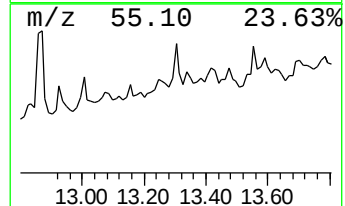
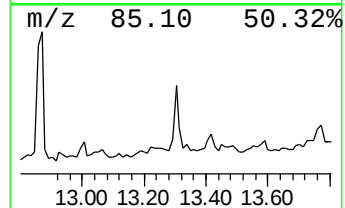
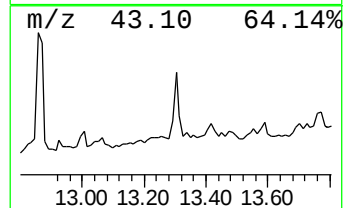
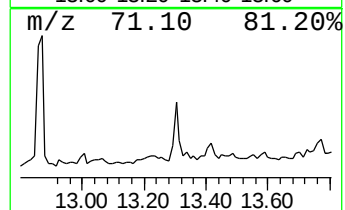
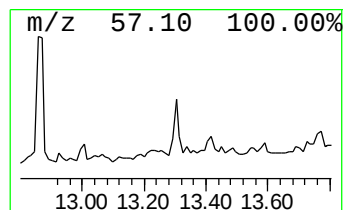
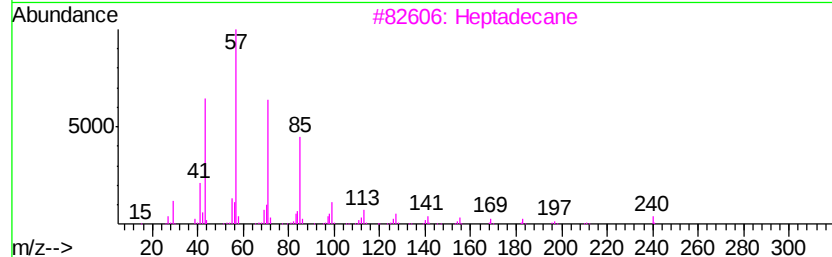
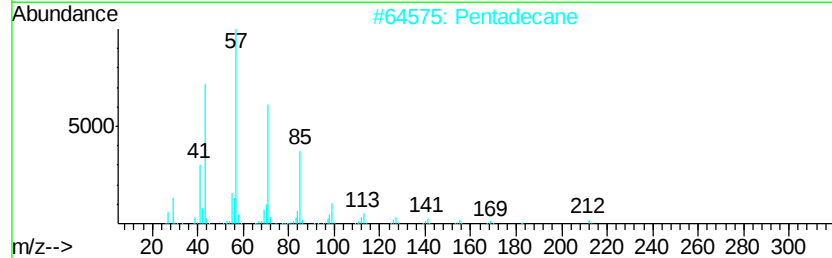
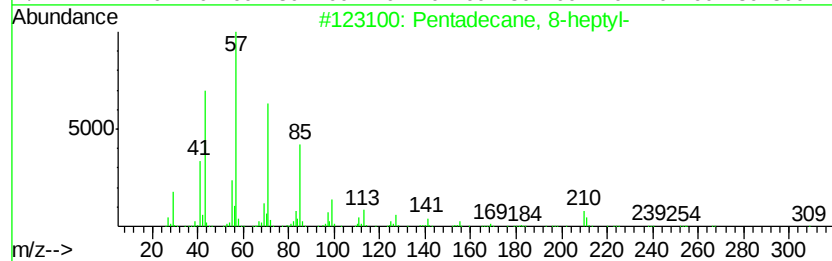
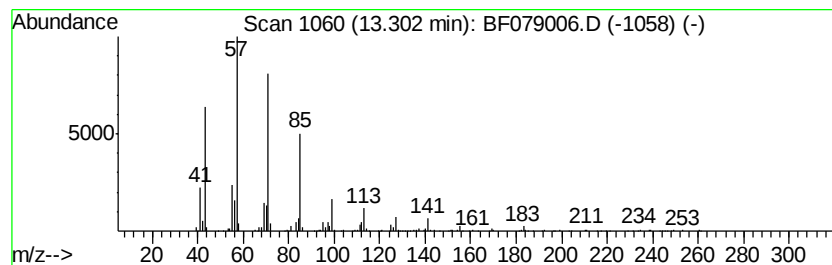
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Pentadecane, 8-heptyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.30	9.54 ng	373800	Phenanthrene-d10	12.95

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecane, 8-heptyl-		310	C22H46	071005-15-7	90
2	Pentadecane		212	C15H32	000629-62-9	90
3	Heptadecane		240	C17H36	000629-78-7	90
4	Octacosane		394	C28H58	000630-02-4	86
5	Tetradecane		198	C14H30	000629-59-4	86



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050715\
Data File : BF079006.D
Acq On : 7 May 2015 20:07
Operator : TP/IZ
Sample : G2084-01 5X
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
1

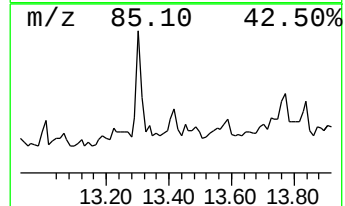
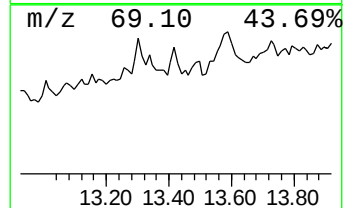
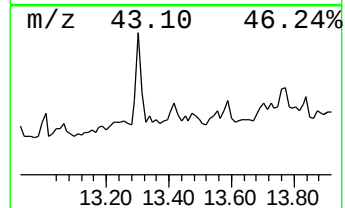
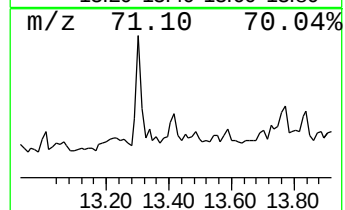
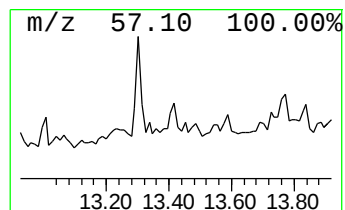
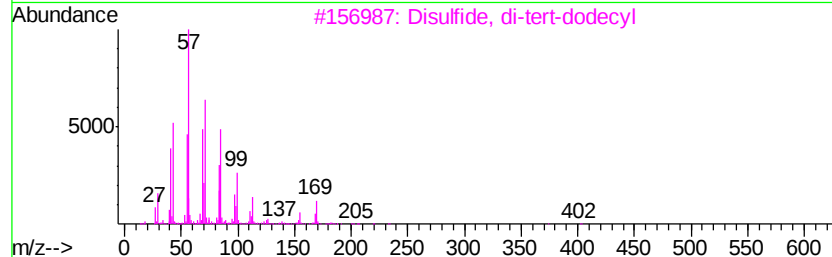
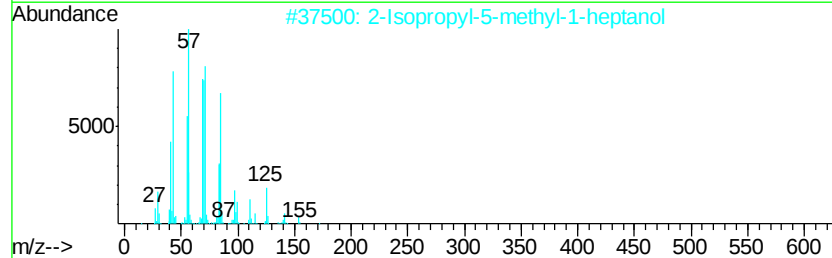
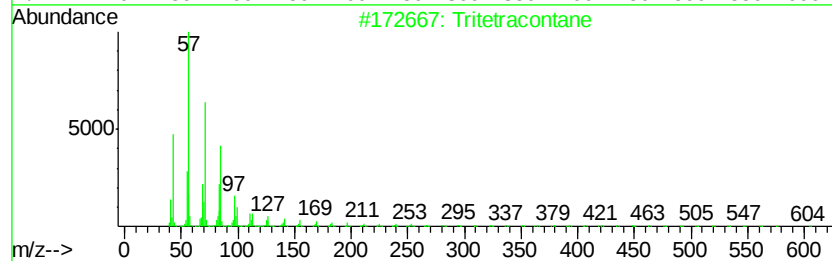
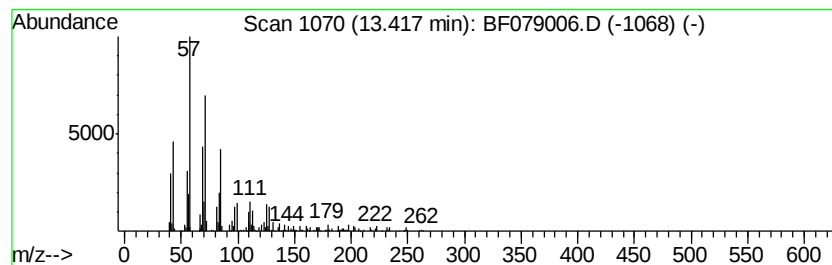
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF043015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 Tritetracontane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.42	3.84 ng	150712	Phenanthrene-d10	12.95

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tritetracontane	605	C43H88	007098-21-7	72
2			2-Isopropyl-5-methyl-1-heptanol	172	C11H24O	091337-07-4	64
3			Disulfide, di-tert-dodecyl	402	C24H50S2	027458-90-8	59
4			Heneicosane	296	C21H44	000629-94-7	58
5			Nonacosane	408	C29H60	000630-03-5	53



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050715\
Data File : BF079006.D
Acq On : 7 May 2015 20:07
Operator : TP/IZ
Sample : G2084-01 5X
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
1

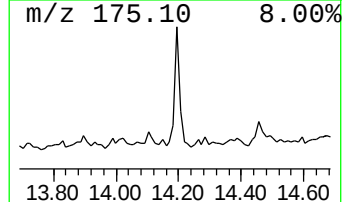
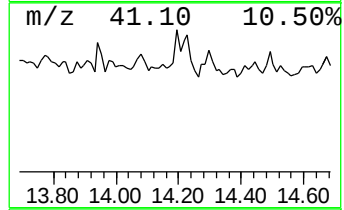
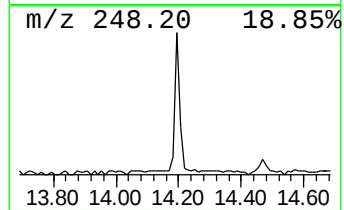
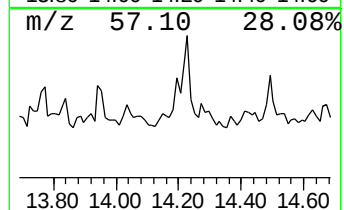
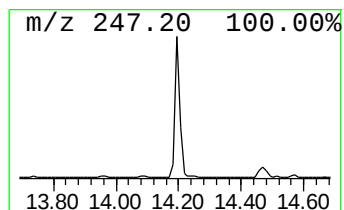
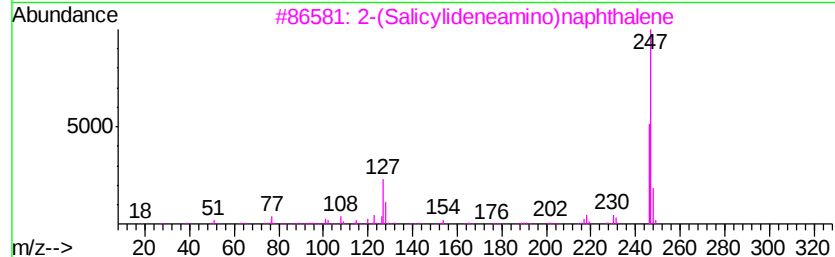
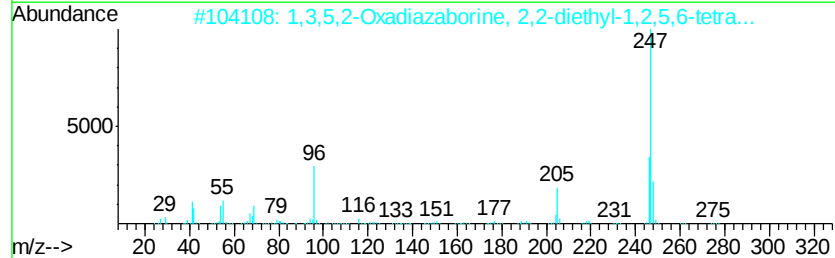
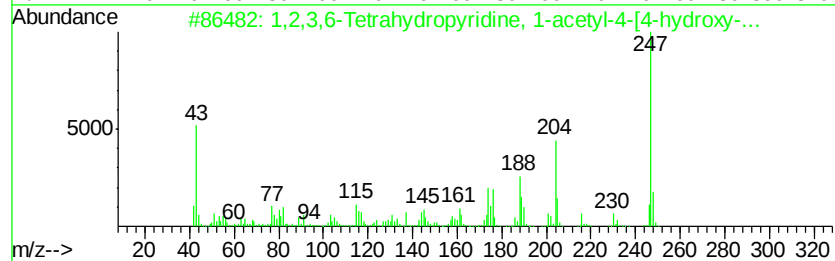
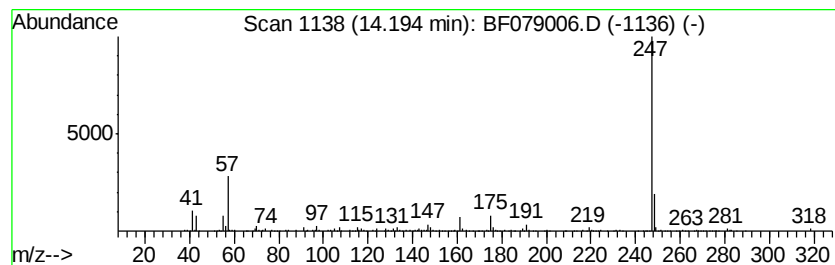
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 1,2,3,6-Tetrahydropyridine,... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.19	6.02 ng	235821	Phenanthrene-d10	12.95

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,2,3,6-Tetrahydropyridine, 1-ac...	247	C14H17N03	094427-34-6	72
2			1,3,5,2-Oxadiazaborine, 2,2-diet...	276	C16H29BN2O	1000160-64-7	64
3			2-(Salicylideneamino)naphthalene	247	C17H13N0	001689-72-1	59
4			9H-Imidazo[1,2-a]benzimidazole, ...	247	C16H13N3	021431-82-3	59
5			ETHYL 4-ANILINO-5-OXO-2,5-DIHYDR...	247	C13H13N04	1000244-19-7	50



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050715\
Data File : BF079006.D
Acq On : 7 May 2015 20:07
Operator : TP/IZ
Sample : G2084-01 5X
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
1

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF043015.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
Propane, 2,2-dime...	1.48	65.4	ng	1178650	1	7.35	360591	20.0
Butane, 2-methoxy...	1.80	8.3	ng	150291	1	7.35	360591	20.0
unknown7.05	7.05	16.6	ng	300173	1	7.35	360591	20.0
unknown11.67	11.67	2.6	ng	79027	3	11.10	612455	20.0
Phenol, 4-(1,1,3,...	11.76	3.1	ng	94793	3	11.10	612455	20.0
Undecane, 3,6-dim...	11.94	7.0	ng	215794	3	11.10	612455	20.0
Phenol, 2,4-bis(1...	12.18	10.3	ng	402645	4	12.95	783993	20.0
Tridecane, 5-propyl-	12.29	13.2	ng	517101	4	12.95	783993	20.0
3-Eicosene, (E)-	12.51	2.6	ng	103153	4	12.95	783993	20.0
Hexadecane	12.87	23.1	ng	903591	4	12.95	783993	20.0
Heptadecane, 2-me...	13.01	2.3	ng	88973	4	12.95	783993	20.0
Pentadecane, 8-he...	13.30	9.5	ng	373800	4	12.95	783993	20.0
Tritetracontane	13.42	3.8	ng	150712	4	12.95	783993	20.0
1,2,3,6-Tetrahydr...	14.19	6.0	ng	235821	4	12.95	783993	20.0