

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF092115\
 Data File : BF081735.D
 Acq On : 22 Sep 2015 5:39
 Operator : UM/IZ
 Sample : G3747-06
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TRENCH-2-2

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-BF090115.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	1.201	8	10	13	rVB	736025	940679	21.94%	5.845%
2	1.281	16	17	29	rVB3	20356	64922	1.51%	0.403%
3	1.441	29	31	37	rBV	134526	359381	8.38%	2.233%
4	1.521	37	38	87	rVB	1157476	4287576	100.00%	26.641%
5	4.424	289	292	310	rVV	190584	548840	12.80%	3.410%
6	5.316	367	370	393	rBV	456875	941953	21.97%	5.853%
7	7.590	566	569	574	rBV	504857	930882	21.71%	5.784%
8	7.682	574	577	591	rVB	554362	1121628	26.16%	6.969%
9	8.162	616	619	625	rBB	123083	221143	5.16%	1.374%
10	8.482	644	647	658	rBB	447122	714853	16.67%	4.442%
11	9.465	729	733	752	rBV	340331	656094	15.30%	4.077%
12	11.054	869	872	879	rBV	150876	277474	6.47%	1.724%
13	13.705	1100	1104	1107	rBV	559686	993509	23.17%	6.173%
14	14.768	1193	1197	1204	rBV	111883	207919	4.85%	1.292%
15	15.191	1230	1234	1237	rBV	164292	294259	6.86%	1.828%
16	17.100	1397	1401	1411	rBV	310806	659635	15.38%	4.099%
17	17.729	1452	1456	1464	rBV	22960	70734	1.65%	0.440%
18	18.700	1537	1541	1544	rBV	171233	306931	7.16%	1.907%
19	19.660	1622	1625	1629	rBV	26857	54377	1.27%	0.338%
20	21.386	1774	1776	1782	rVB	22467	46117	1.08%	0.287%
21	22.072	1833	1836	1838	rBV	732783	1005805	23.46%	6.250%
22	23.306	1942	1944	1946	rBV	71958	134426	3.14%	0.835%
23	23.352	1946	1948	1950	rVB	180485	259268	6.05%	1.611%
24	24.004	2003	2005	2009	rBV	47765	55341	1.29%	0.344%
25	24.598	2055	2057	2064	rBV2	56014	97824	2.28%	0.608%
26	24.781	2072	2073	2080	rVB	127111	189740	4.43%	1.179%
27	25.181	2105	2108	2110	rBV	66906	86147	2.01%	0.535%
28	25.821	2158	2164	2168	rBV4	14003	47291	1.10%	0.294%
29	25.924	2169	2173	2176	rVV2	23276	59177	1.38%	0.368%
30	26.015	2179	2181	2188	rVB4	30884	73259	1.71%	0.455%
31	26.198	2191	2197	2201	rBV3	38405	101568	2.37%	0.631%
32	26.472	2217	2221	2226	rBV3	17226	44629	1.04%	0.277%
33	26.884	2253	2257	2266	rVB3	13382	44910	1.05%	0.279%
34	27.535	2310	2314	2322	rVB4	22273	57357	1.34%	0.356%

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ALS Vial : 23 Sample Multiplier: 1

Instrument :
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ClientSampleId :
TRENCH-2-2

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-BF090115.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	27.798	2332	2337	2343	rVB	52723	138313	3.23%	0.859%
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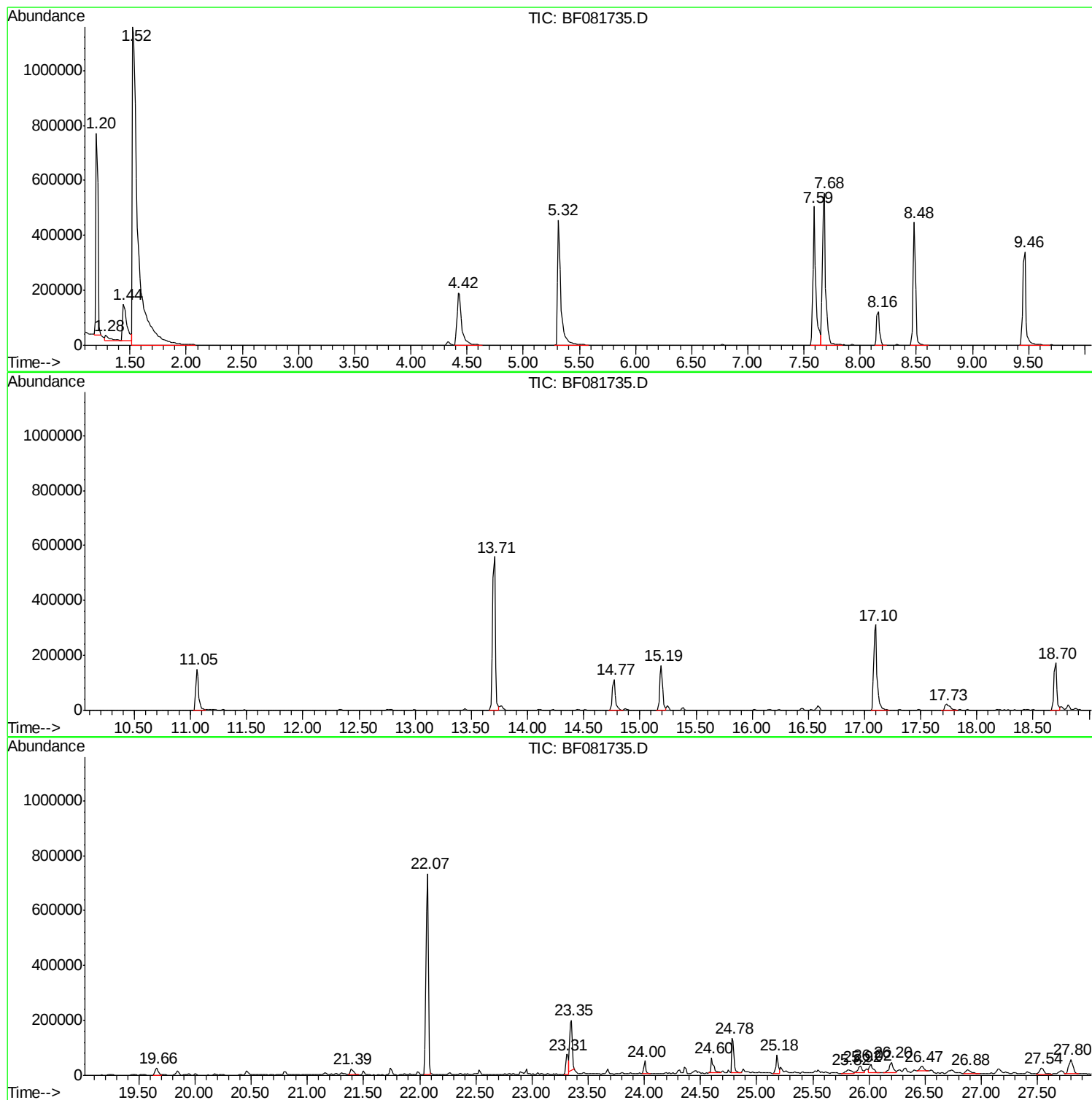
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF090115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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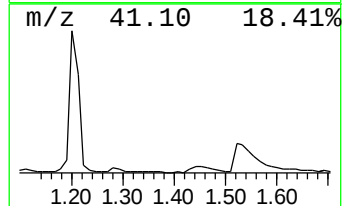
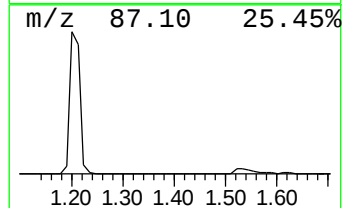
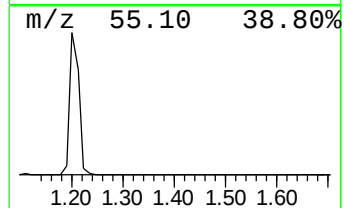
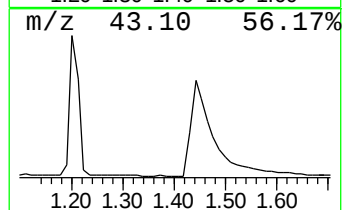
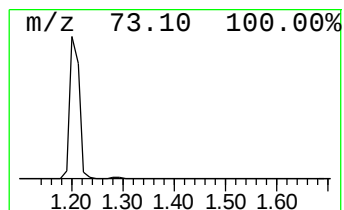
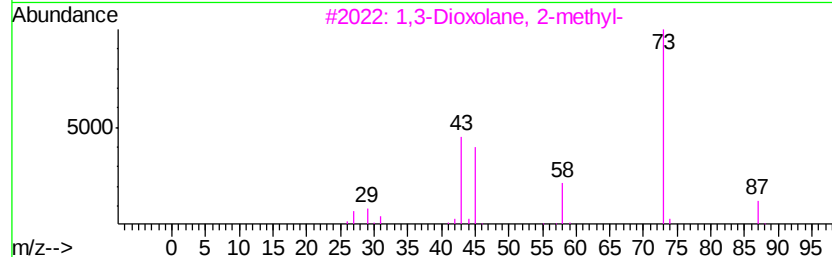
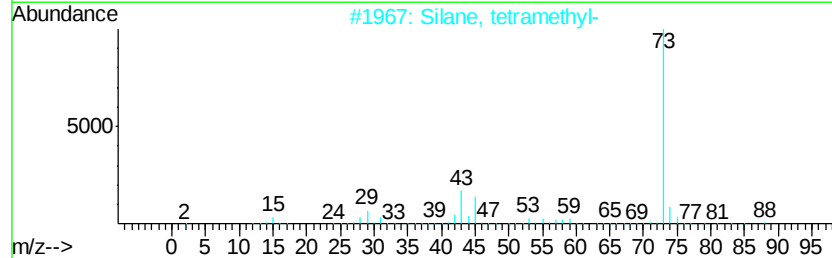
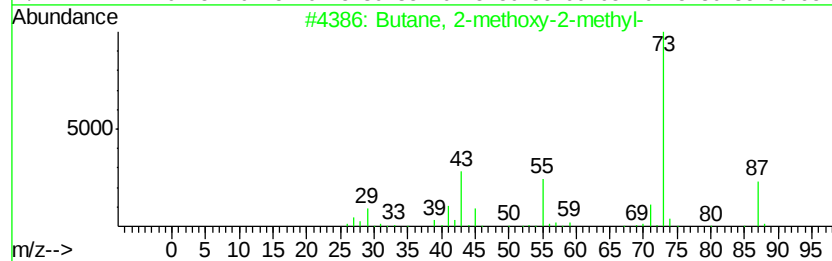
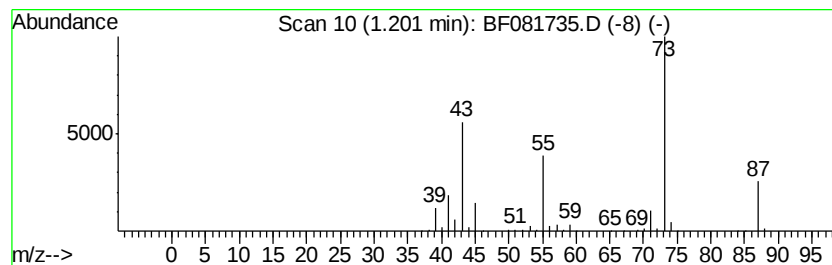
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.	
1.20	85.07 ng	940679	1,4-Dichlorobenzene-d4	8.16	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2	Silane, tetramethyl-	88	C4H12Si	000075-76-3	12
3	1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9
4	sec-Butyl nitrite	103	C4H9NO2	000924-43-6	9
5	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	9



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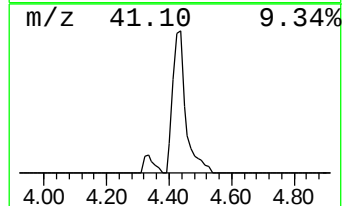
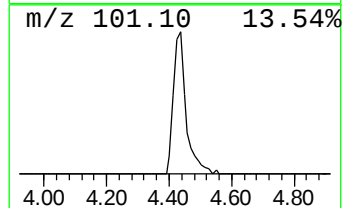
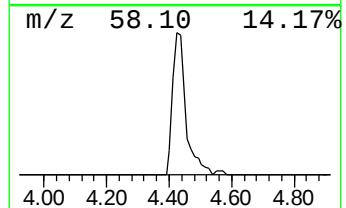
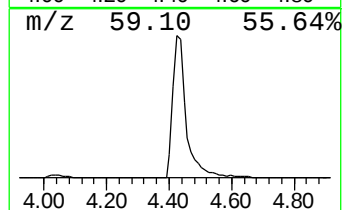
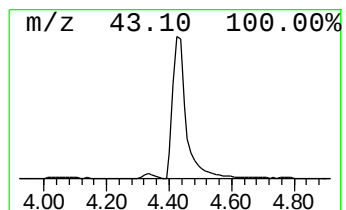
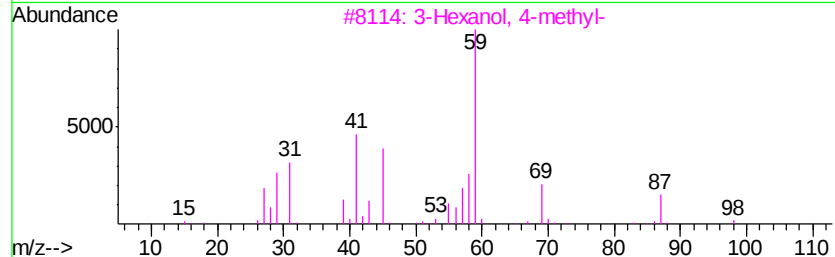
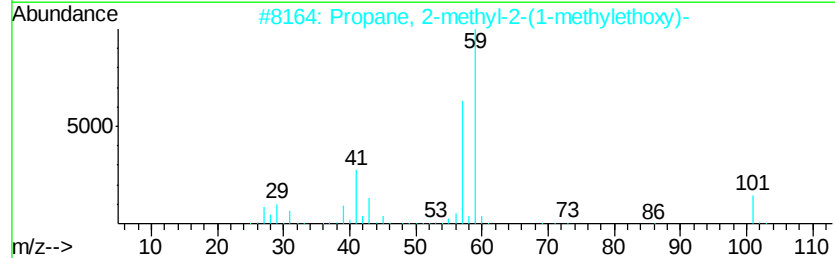
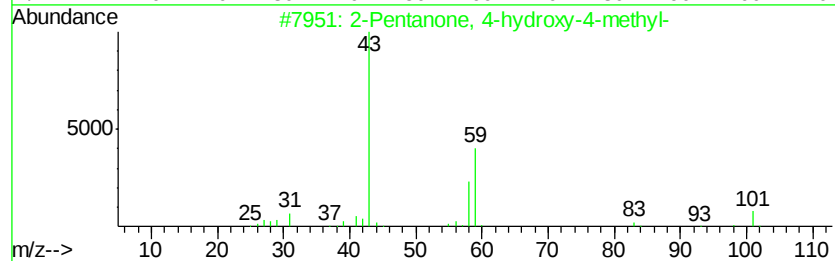
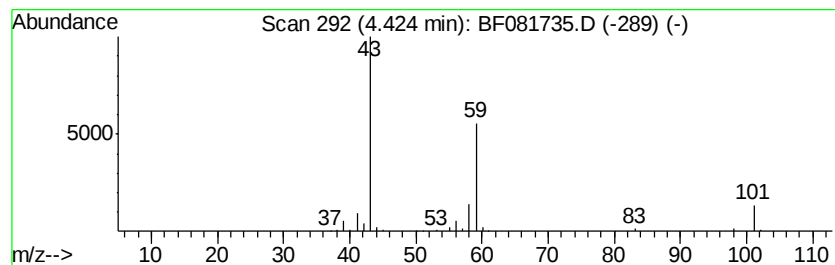
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.42	49.64 ng	548840	1,4-Dichlorobenzene-d4	8.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2		Propane, 2-methyl-2-(1-methylethoxy)-	116	C7H16O	017348-59-3	42
3		3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
4		Acetic acid, cyano-, 1,1-dimethyl-	141	C7H11NO2	001116-98-9	25
5		Acetic acid, 1,1-dimethylethyl ester	116	C6H12O2	000540-88-5	25



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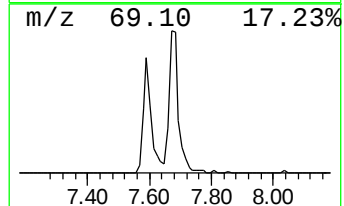
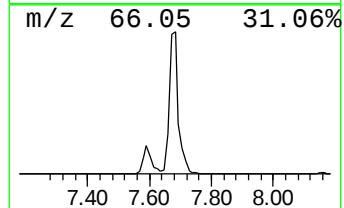
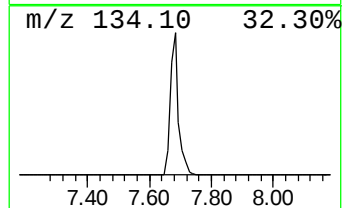
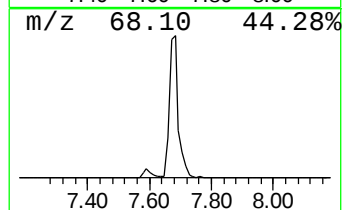
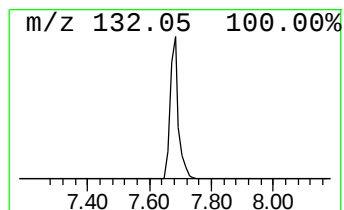
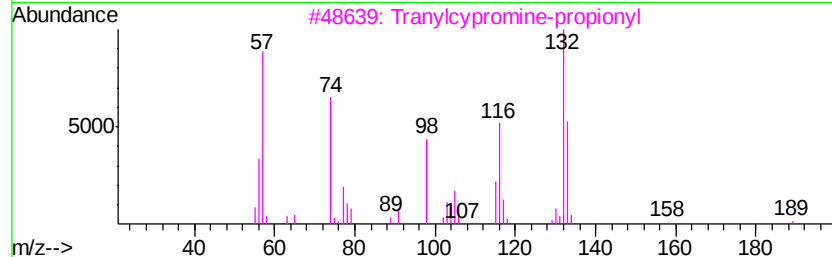
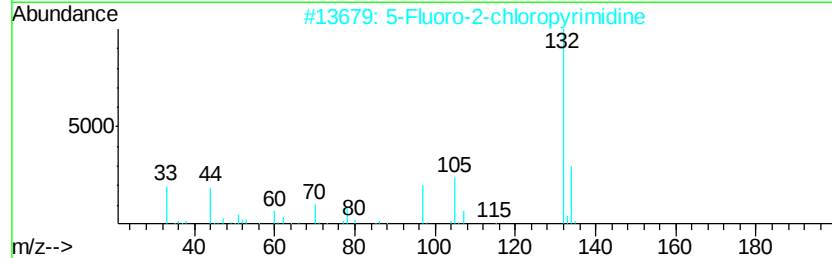
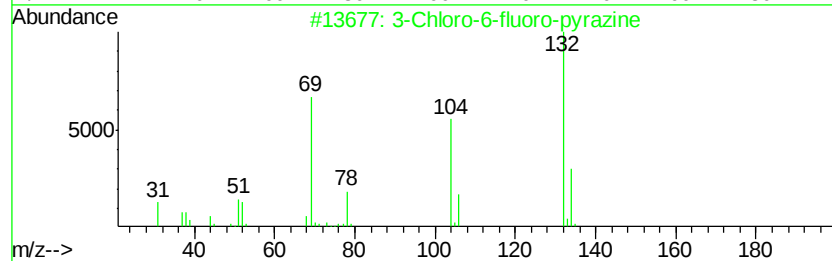
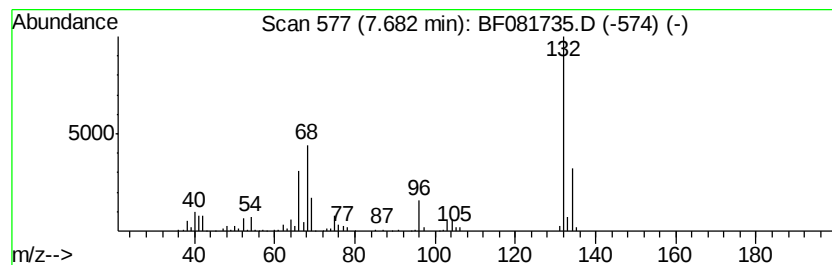
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 unknown7.68 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.68	101.44 ng	1121630	1,4-Dichlorobenzene-d4	8.16

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	12
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	10
4		(5-METHYL-2-PYRIDYL)ACETONITRILE	132	C8H8N2	1000241-93-9	10
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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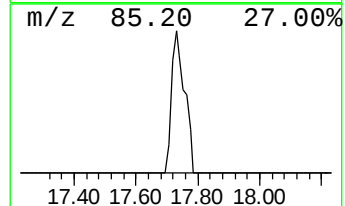
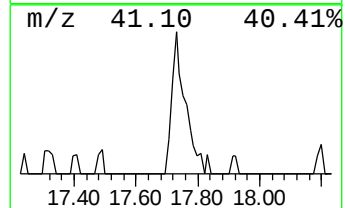
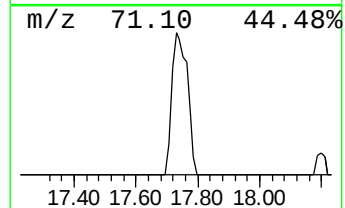
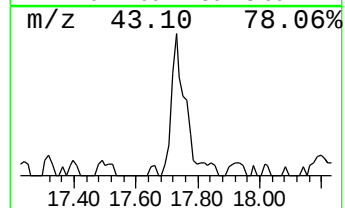
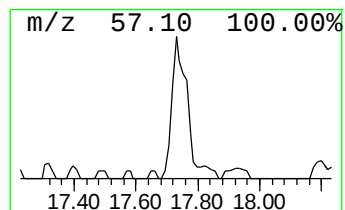
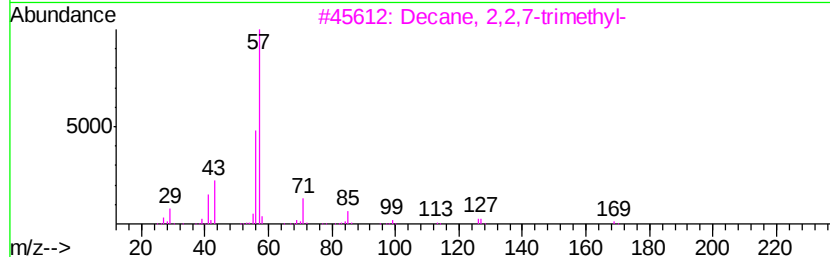
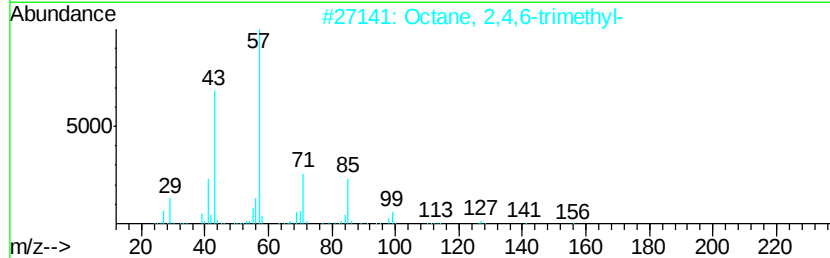
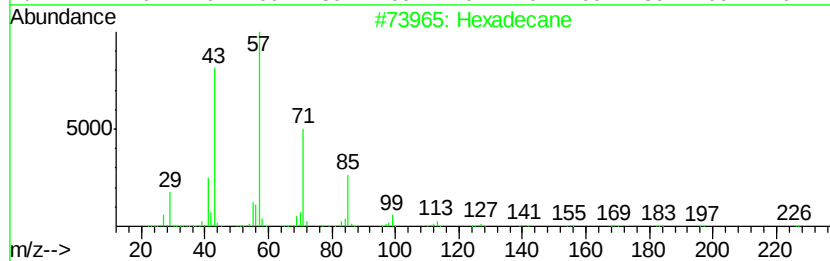
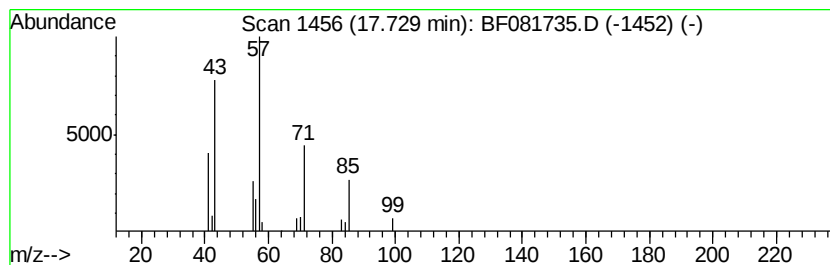
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Hexadecane Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.73	4.61 ng	70734	Phenanthrene-d10	18.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecane	226	C16H34	000544-76-3	83
2		Octane, 2,4,6-trimethyl-	156	C11H24	062016-37-9	64
3		Decane, 2,2,7-trimethyl-	184	C13H28	062237-99-4	59
4		Decane, 2,5,9-trimethyl-	184	C13H28	062108-22-9	59
5		Tridecane	184	C13H28	000629-50-5	56



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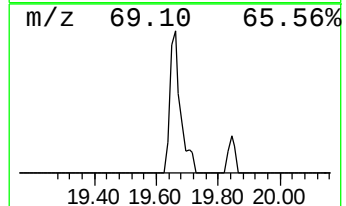
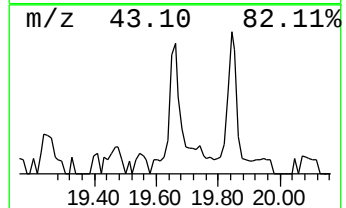
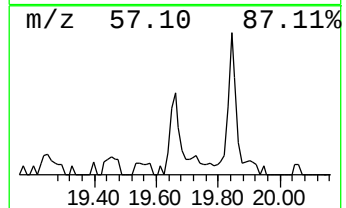
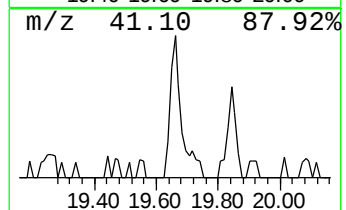
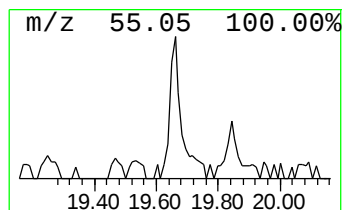
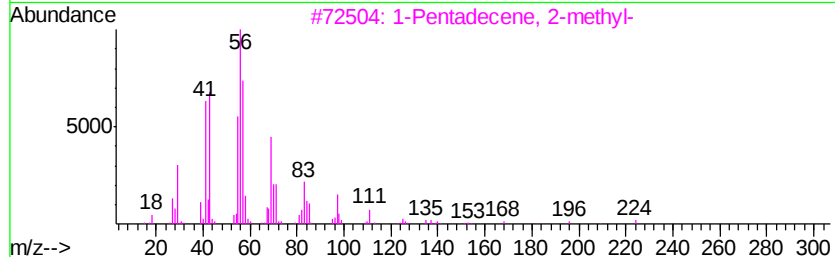
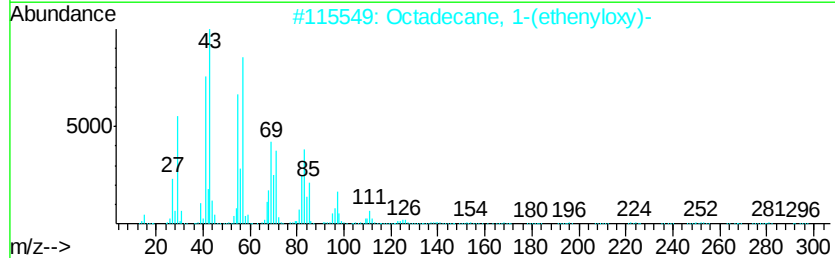
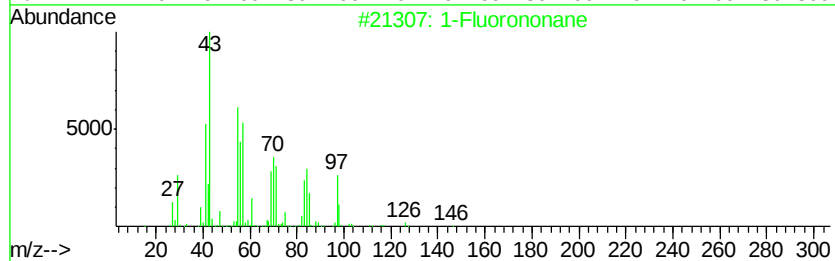
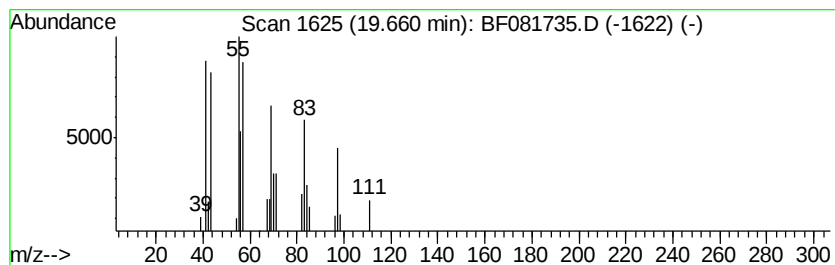
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 9 unknown19.66 Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.66	3.54 ng	54377	Phenanthrene-d10	18.70

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Fluorononane		146	C9H19F	000463-18-3	47
2	Octadecane, 1-(ethenvloxy)-		296	C20H40O	000930-02-9	43
3	1-Pentadecene, 2-methyl-		224	C16H32	029833-69-0	38
4	Silane, trichlorodocosyl-		442	C22H45Cl3Si	007325-84-0	37
5	1-Heptanol, 6-methyl-		130	C8H18O	001653-40-3	27



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ALS Vial : 23 Sample Multiplier: 1

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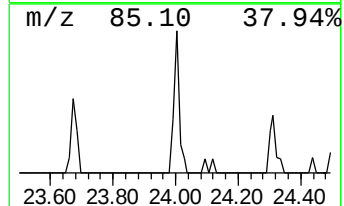
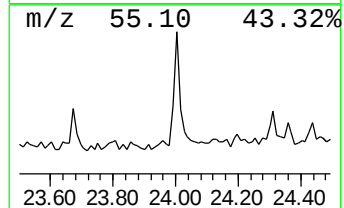
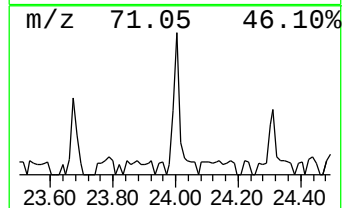
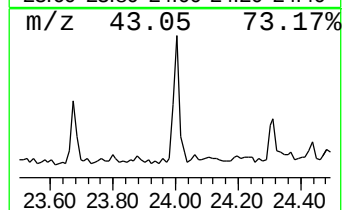
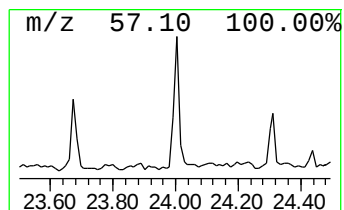
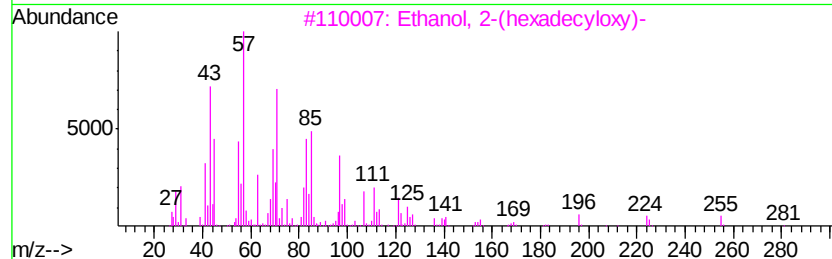
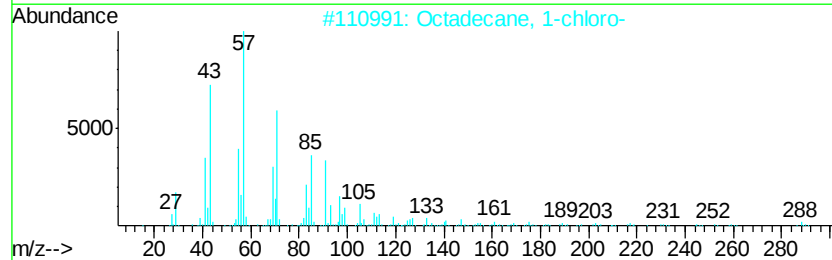
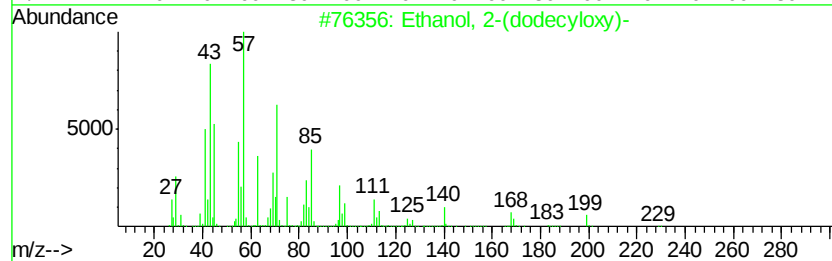
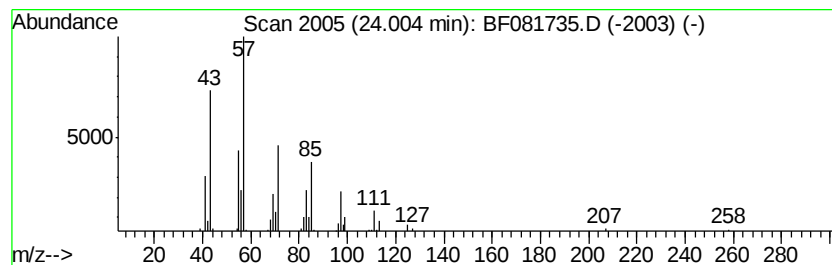
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Ethanol, 2-(dodecyloxy)- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.00	4.27 ng	55341	Chrysene-d12	23.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Ethanol, 2-(dodecyloxy)-	230	C14H30O2	004536-30-5	86
2		Octadecane, 1-chloro-	288	C18H37Cl	003386-33-2	72
3		Ethanol, 2-(hexadecyloxy)-	286	C18H38O2	002136-71-2	72
4		2-Ethyl-1-dodecanol	214	C14H30O	019780-33-7	72
5		Tetradecane	198	C14H30	000629-59-4	52



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF092115\
Data File : BF081735.D
Acq On : 22 Sep 2015 5:39
Operator : UM/IZ
Sample : G3747-06
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TRENCH-2-2

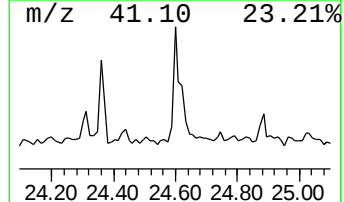
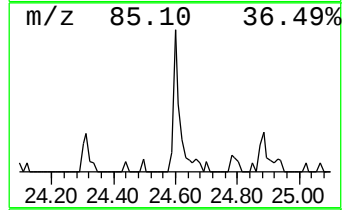
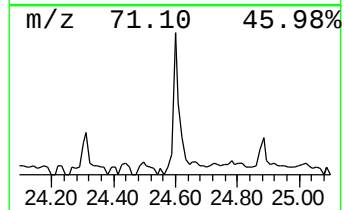
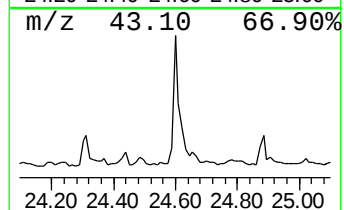
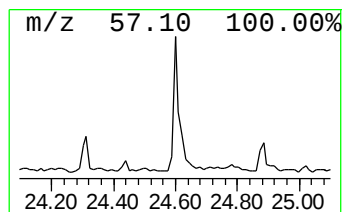
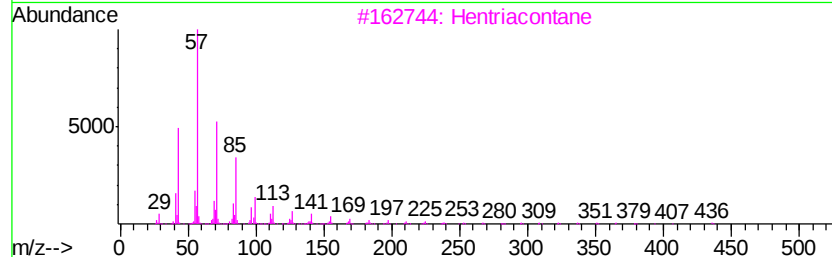
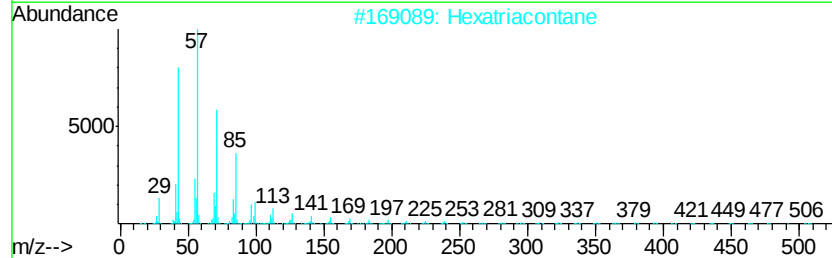
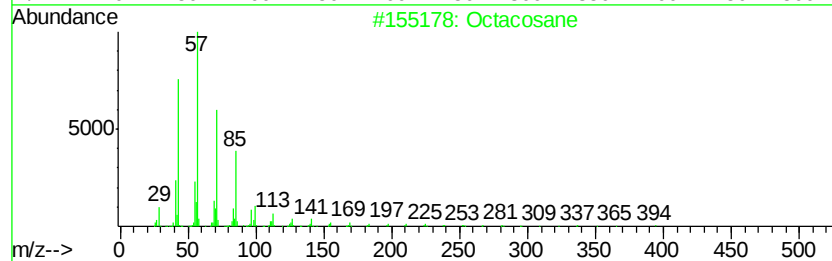
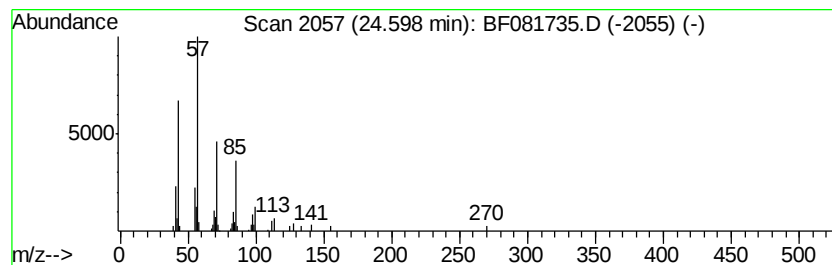
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Octacosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.60	10.31 ng	97824	Perylene-d12	24.79

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octacosane	394	C28H58	000630-02-4	91
2		Hexatriacontane	507	C36H74	000630-06-8	91
3		Hentriacontane	437	C31H64	000630-04-6	90
4		Heptacosane	380	C27H56	000593-49-7	87
5		Eicosane	282	C20H42	000112-95-8	87



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF092115\
Data File : BF081735.D
Acq On : 22 Sep 2015 5:39
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Misc :
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Instrument :
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ClientSampleId :
TRENCH-2-2

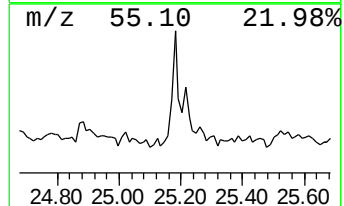
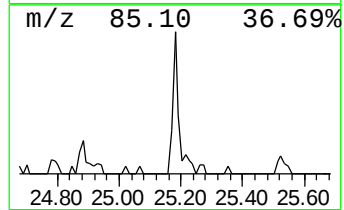
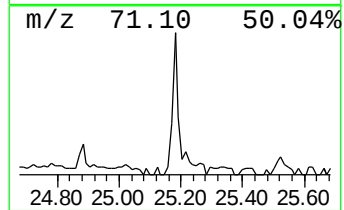
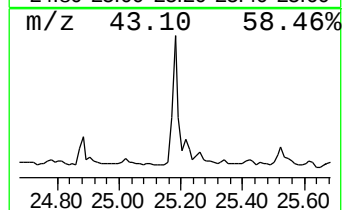
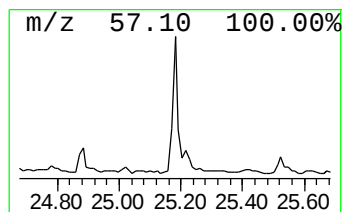
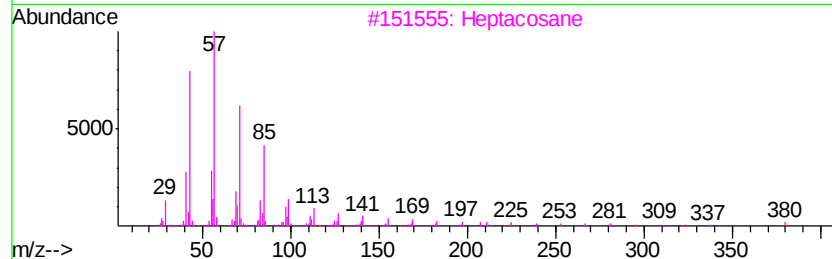
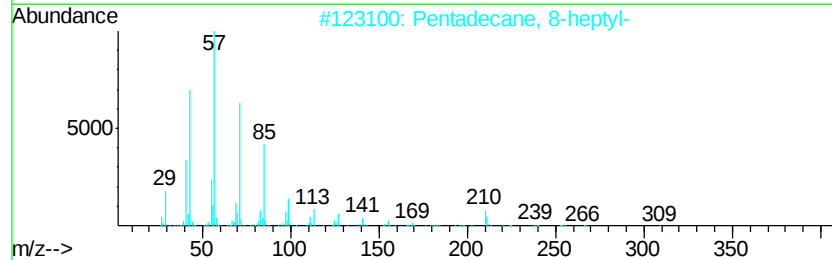
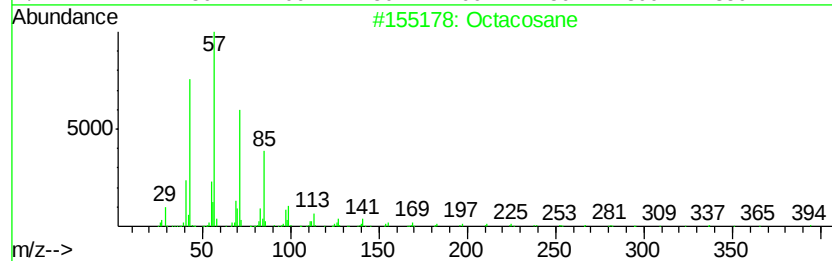
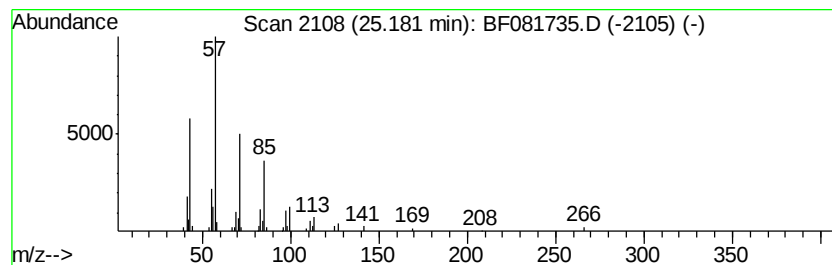
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Pentadecane, 8-heptyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.18	9.08 ng	86147	Perylene-d12	24.79

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octacosane	394	C28H58	000630-02-4	90
2		Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	90
3		Heptacosane	380	C27H56	000593-49-7	87
4		Tetracosane	338	C24H50	000646-31-1	87
5		Tetratetracontane	619	C44H90	007098-22-8	86



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF092115\
Data File : BF081735.D
Acq On : 22 Sep 2015 5:39
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Sample : G3747-06
Misc :
ALS Vial : 23 Sample Multiplier: 1

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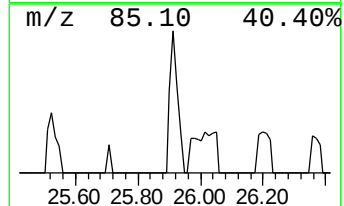
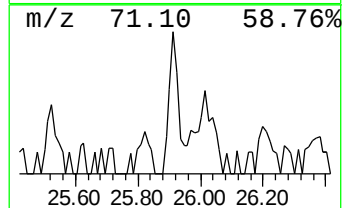
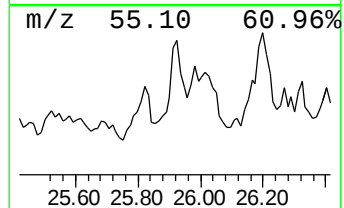
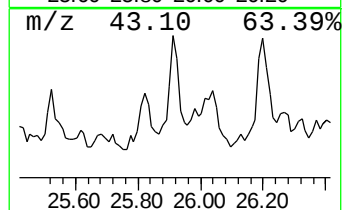
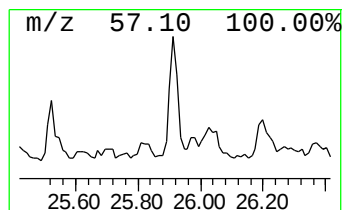
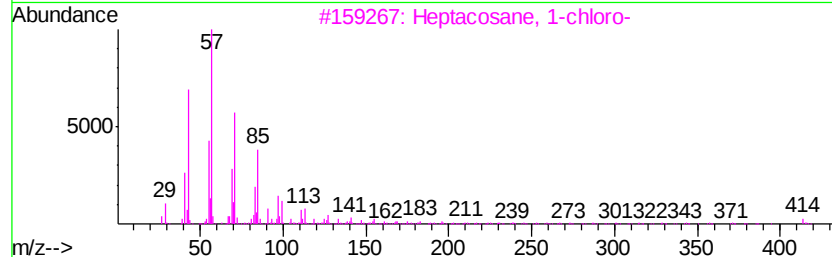
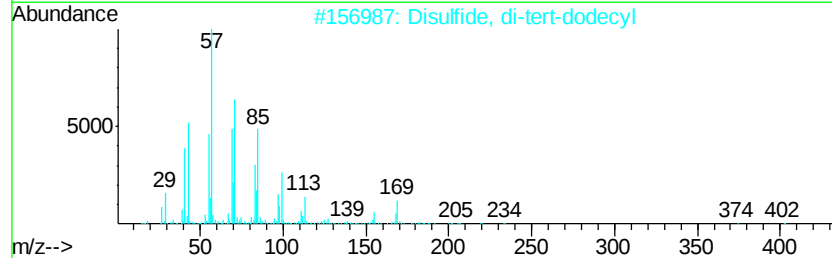
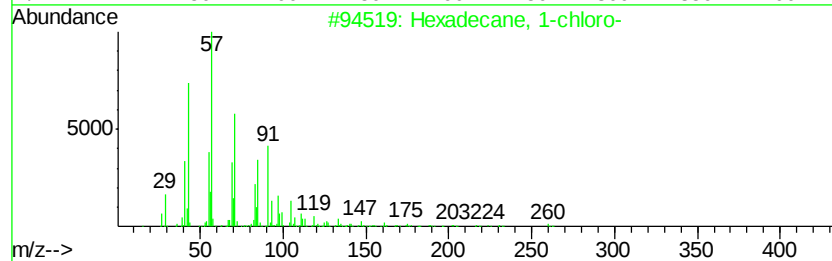
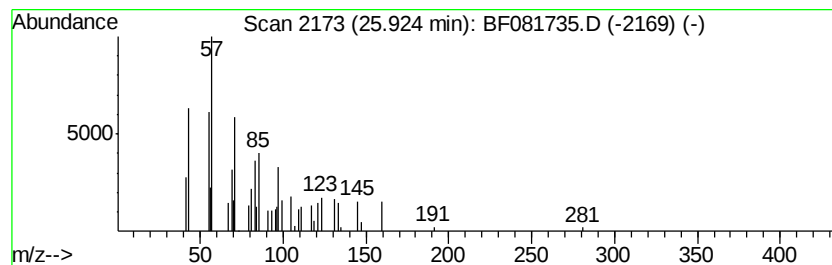
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 unknown25.92 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.92	6.24 ng	59177	Perylene-d12	24.79

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecane, 1-chloro-	260	C16H33Cl	004860-03-1	43
2		Disulfide, di-tert-dodecyl	402	C24H50S2	027458-90-8	38
3		Heptacosane, 1-chloro-	414	C27H55Cl	062016-79-9	38
4		Nonadecane	268	C19H40	000629-92-5	38
5		Tetradecane	198	C14H30	000629-59-4	35



Data Path : Z:\HPCHEM1\BNA F\DATA\BF092115\
Data File : BF081735.D
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ClientSampleId :
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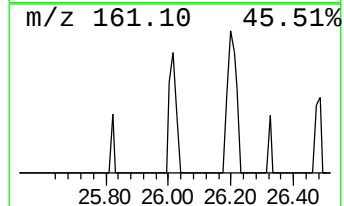
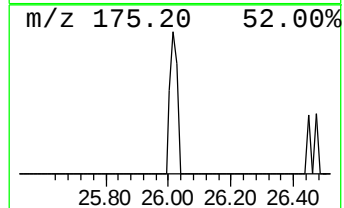
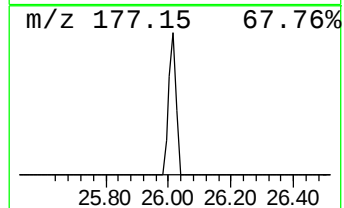
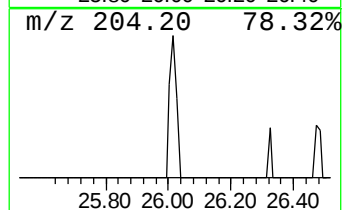
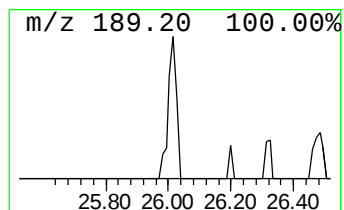
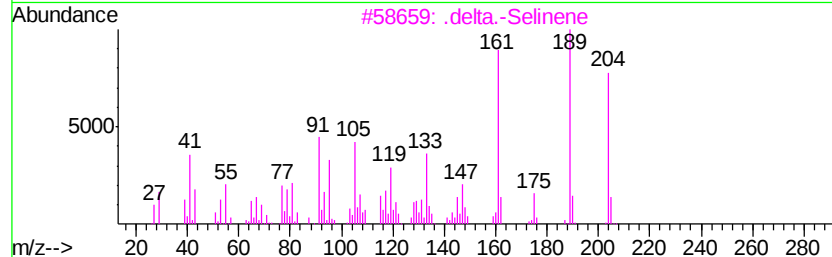
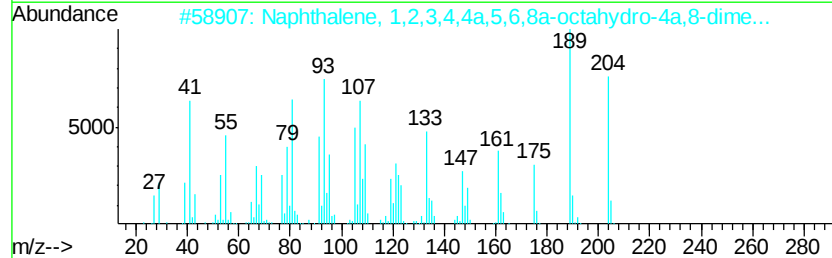
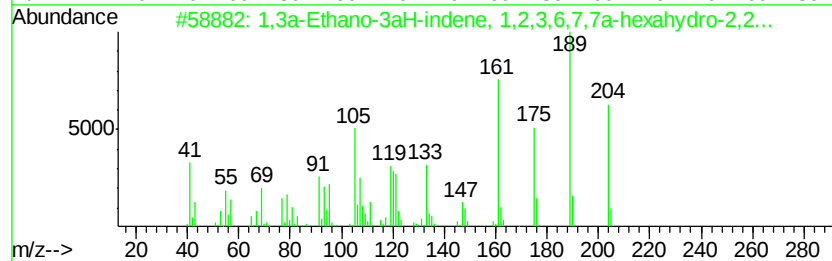
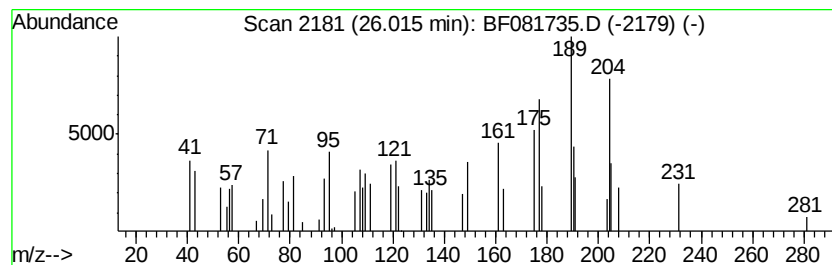
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 1,3a-Ethano-3aH-indene, 1,2... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.02	7.72 ng	73259	Perylene-d12	24.79

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,3a-Ethano-3aH-indene, 1,2,3,6,...	204	C15H24	004545-68-0	70
2		Naphthalene, 1,2,3,4,4a,5,6,8a-o...	204	C15H24	000473-13-2	50
3		.delta.-Selinene	204	C15H24	028624-23-9	50
4		1H-Cyclopropa[alnaphthalene, 1a,...	204	C15H24	000489-29-2	46
5		1H-Cyclopropa[a]naphthalene, 1a,...	204	C15H24	1000147-32-8	43



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF092115\
Data File : BF081735.D
Acq On : 22 Sep 2015 5:39
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ClientSampleId :
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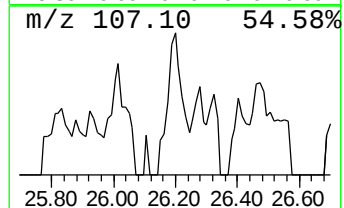
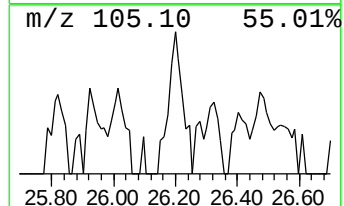
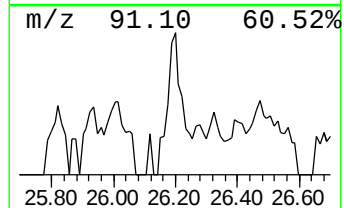
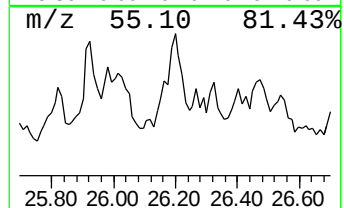
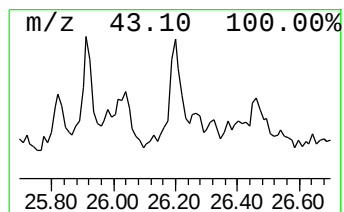
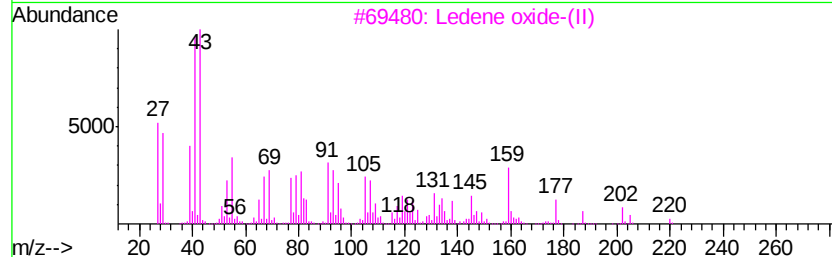
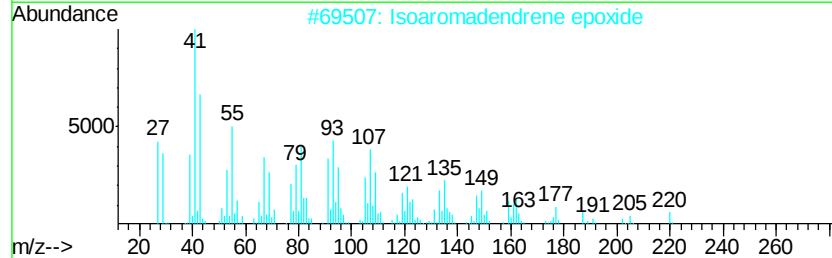
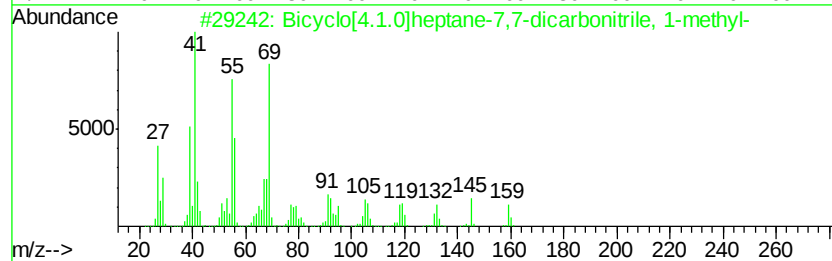
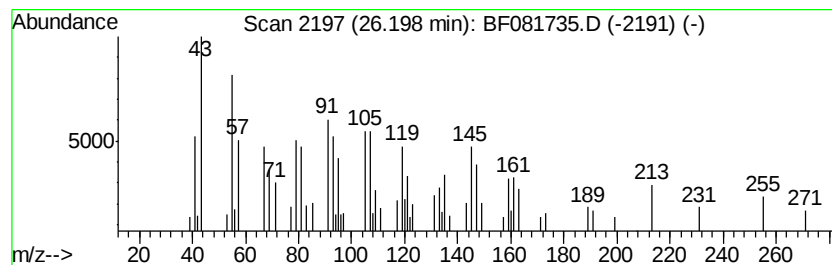
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 unknown26.20 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.20	10.71 ng	101568	Perylene-d12	24.79

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Bicyclo[4.1.0]heptane-7,7-dicarb...	160	C10H12N2	074764-53-7	35
2		Isoaromadendrene epoxide	220	C15H24O	1000159-36-6	27
3		Ledene oxide-(II)	220	C15H24O	1000159-36-7	22
4		2H-Pyran, 2-(7-heptadecynyl)oxv)t...	336	C22H40O2	056599-50-9	12
5		Rishitin	222	C14H22O2	018178-54-6	10



Data Path : Z:\HPCHEM1\BNA F\DATA\BF092115\
Data File : BF081735.D
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Misc :
ALS Vial : 23 Sample Multiplier: 1

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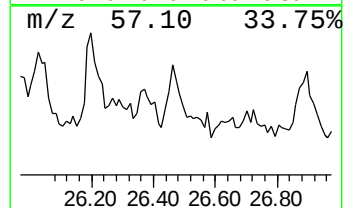
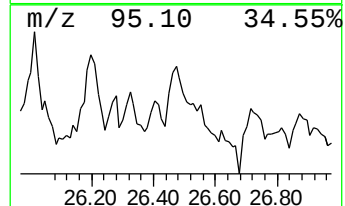
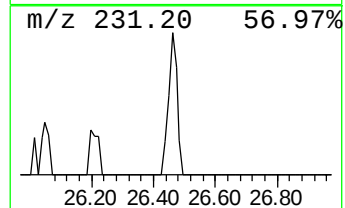
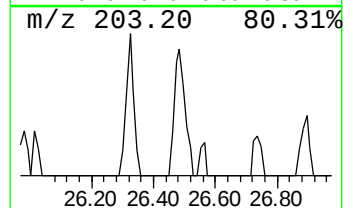
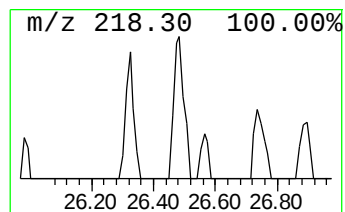
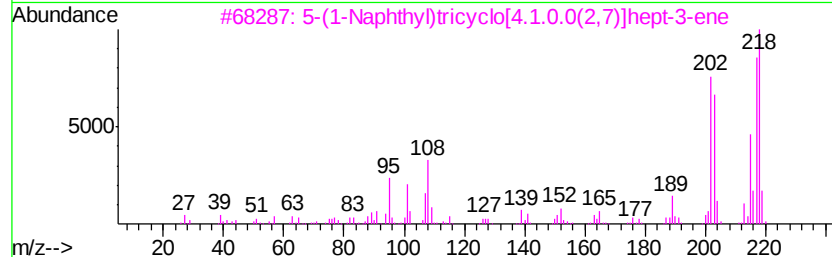
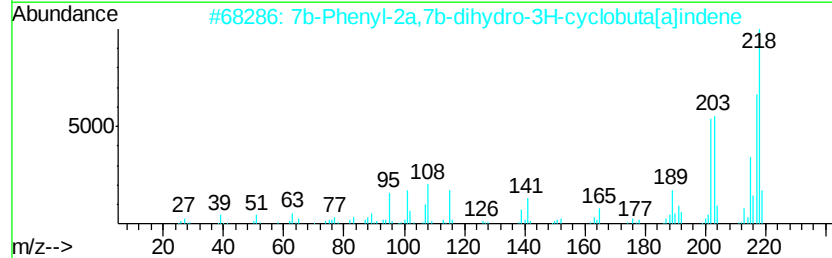
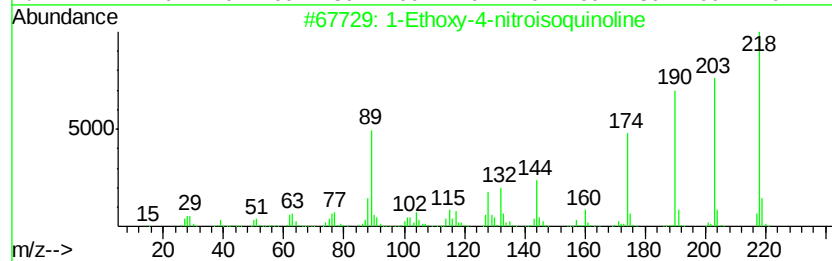
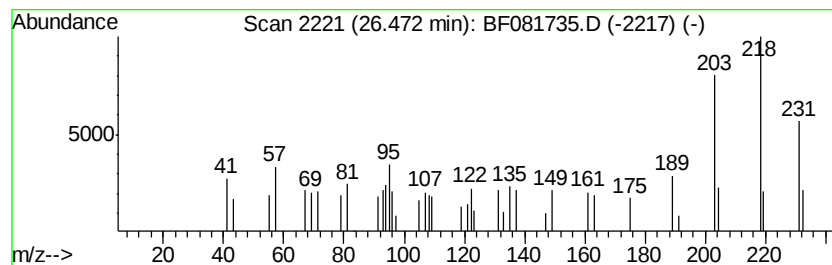
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 unknown26.47 Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.47	4.70 ng	44629	Perylene-d12	24.79

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Ethoxy-4-nitroisoquinoline	218	C11H10N2O3	103863-04-3	43
2			7b-Phenyl-2a,7b-dihydro-3H-cyclo...	218	C17H14	138089-70-0	43
3			5-(1-Naphthyl)tricyclo[4.1.0.0(2...	218	C17H14	107729-05-5	43
4			3-Methoxy-6,7,8,9-tetrahydro-dib...	218	C13H14O3	1000186-57-9	43
5			2H-Cyclopropa[a]naphthalen-2-one...	218	C15H22O	006831-17-0	38



Data Path : Z:\HPCHEM1\BNA F\DATA\BF092115\
Data File : BF081735.D
Acq On : 22 Sep 2015 5:39
Operator : UM/IZ
Sample : G3747-06
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TRENCH-2-2

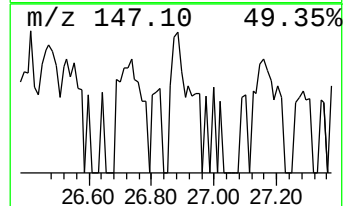
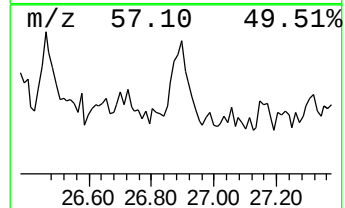
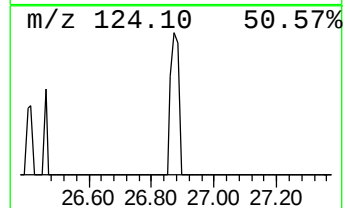
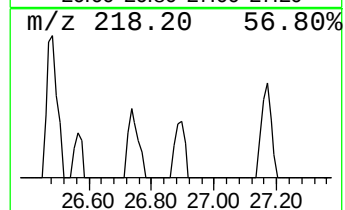
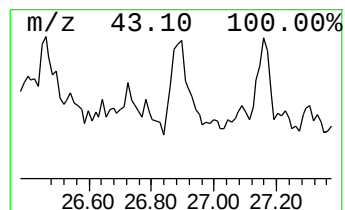
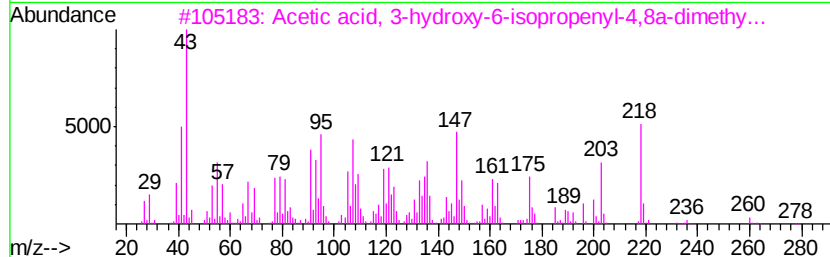
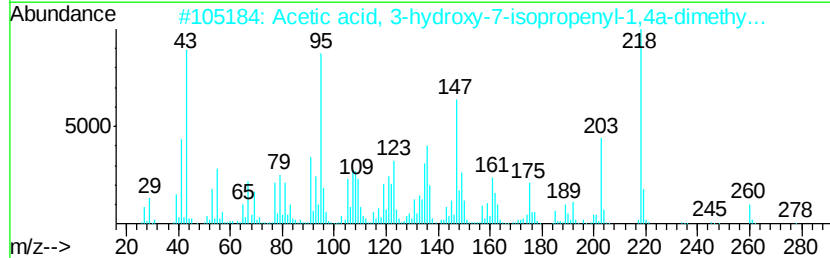
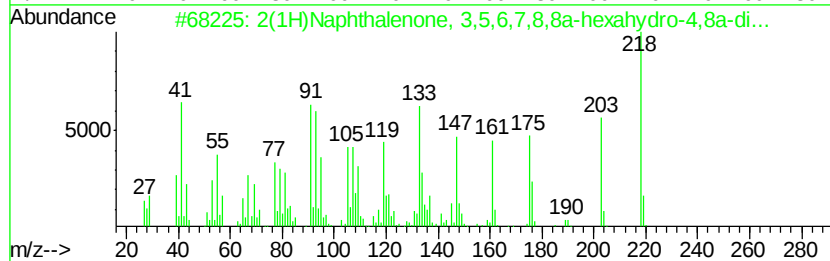
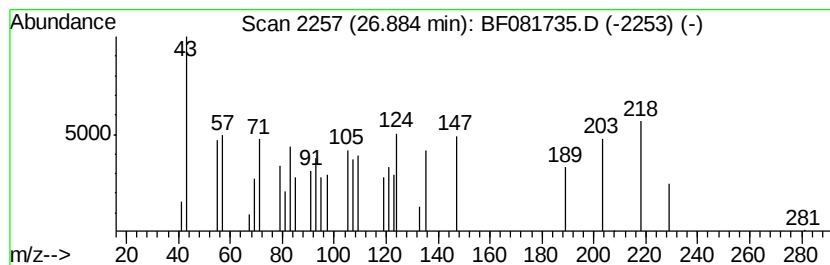
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF090115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 unknown26.88 Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.88	4.73 ng	44910	Perylene-d12	24.79

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2(1H)Naphthalenone, 3,5,6,7,8,8a...	218	C15H22O	1000188-66-5	27
2		Acetic acid, 3-hydroxy-7-isoprop...	278	C17H26O3	1000187-37-6	25
3		Acetic acid, 3-hydroxy-6-isoprop...	278	C17H26O3	1000185-44-8	20
4		5H-3,5a-Epoxy naphth[2,1-c]oxepin...	278	C18H30O2	001153-34-0	12
5		Ledol	222	C15H26O	000577-27-5	10



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF092115\
Data File : BF081735.D
Acq On : 22 Sep 2015 5:39
Operator : UM/IZ
Sample : G3747-06
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TRENCH-2-2

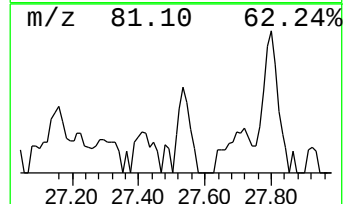
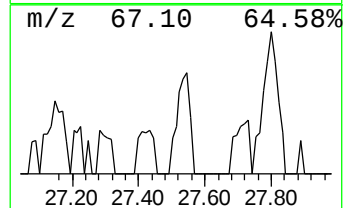
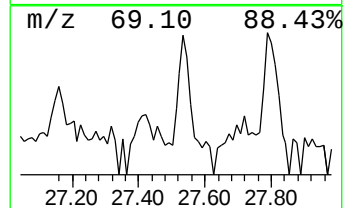
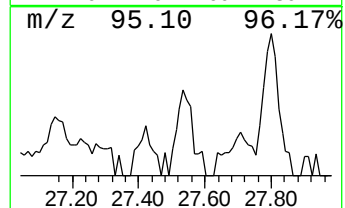
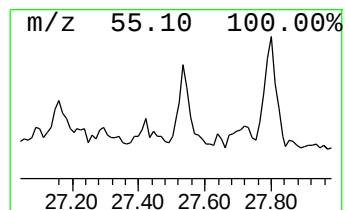
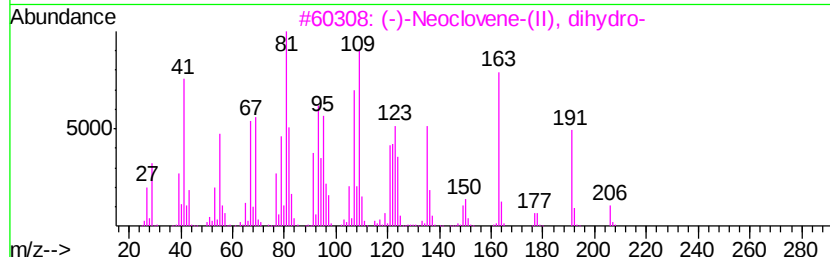
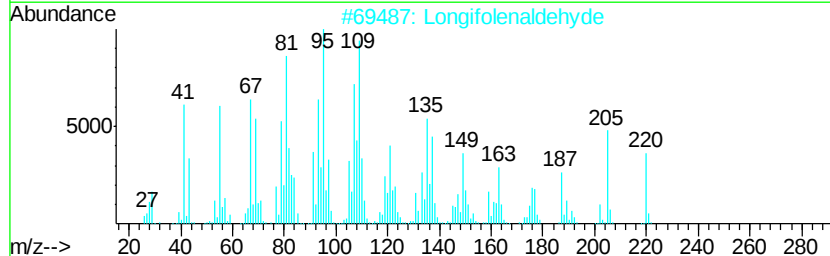
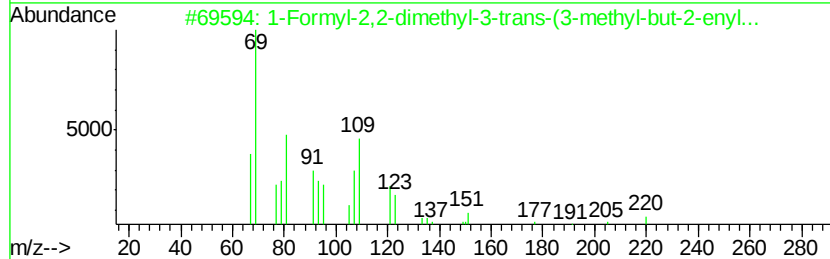
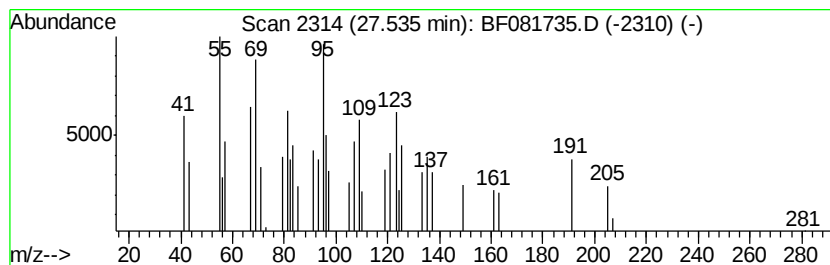
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 unknown27.54 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
27.54	6.05 ng	57357	Perylene-d12	24.79

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Formyl-2,2-dimethyl-3-trans-(3...	220	C15H24O	1000144-09-7	35
2			Longifolenaldehyde	220	C15H24O	019890-84-7	30
3			(-)-Neoclovene-(II), dihydro-	206	C15H26	1000152-83-6	27
4			2,4,4-Trimethyl-3-hydroxymethyl-...	222	C15H26O	1000144-10-5	27
5			(R)-(-)-14-Methyl-8-hexadecyn-1-ol	252	C17H32O	064566-18-3	27



Data Path : Z:\HPCHEM1\BNA F\DATA\BF092115\
Data File : BF081735.D
Acq On : 22 Sep 2015 5:39
Operator : UM/IZ
Sample : G3747-06
Misc :
ALS Vial : 23 Sample Multiplier: 1

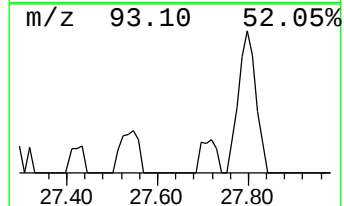
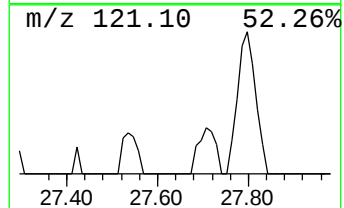
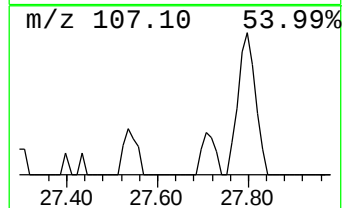
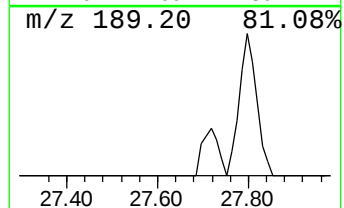
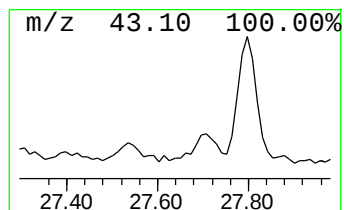
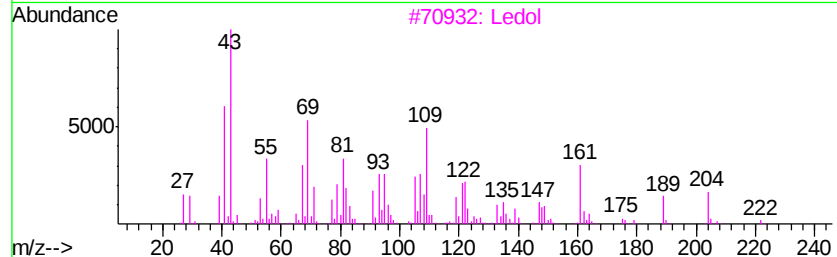
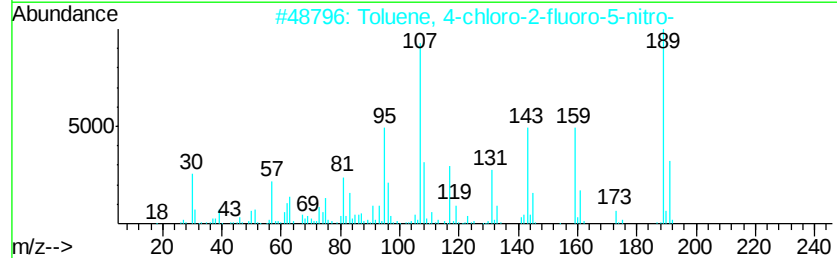
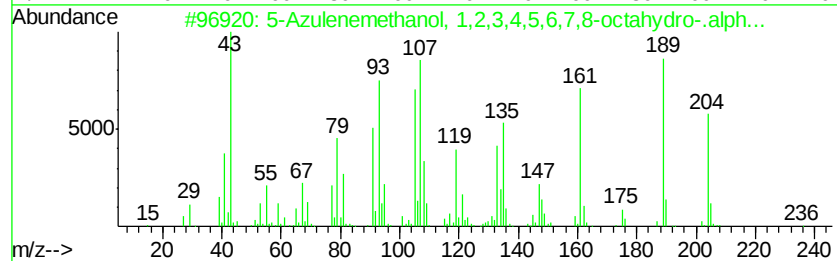
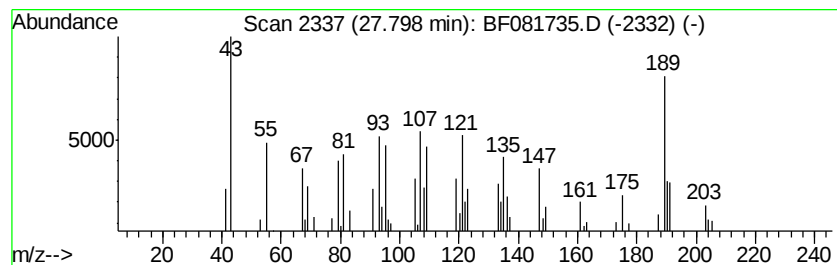
Instrument :
BNA_F
ClientSampleId :
TRENCH-2-2

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF090115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 20 unknown27.80 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.		
27.80	14.58 ng	138313	Perylene-d12	24.79		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Azulenemethanol,	1,2,3,4,5,6,7...	264	C17H28O2	000134-28-1	43
2	Toluene, 4-chloro-2-fluoro-5-nitro-		189	C7H5ClFN02	018349-11-6	43
3	Ledol		222	C15H26O	000577-27-5	27
4	Ethanone, 1-[4-(trifluoromethoxyv...		204	C9H7F3O2	085013-98-5	27
5	Cycloheptane, 4-methylene-1-meth...		204	C15H24	1000159-38-5	25



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF092115\
 Data File : BF081735.D
 Acq On : 22 Sep 2015 5:39
 Operator : UM/IZ
 Sample : G3747-06
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TRENCH-2-2

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF090115.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
Butane, 2-methoxy...	1.20	85.1	ng	940679	1	8.16	221143	20.0
2-Pentanone, 4-hy...	4.42	49.6	ng	548840	1	8.16	221143	20.0
unknown7.68	7.68	101.4	ng	1121630	1	8.16	221143	20.0
Hexadecane	17.73	4.6	ng	70734	4	18.70	306931	20.0
unknown19.66	19.66	3.5	ng	54377	4	18.70	306931	20.0
Ethanol, 2-(dodec...	24.00	4.3	ng	55341	5	23.35	259268	20.0
Octacosane	24.60	10.3	ng	97824	6	24.79	189740	20.0
Pentadecane, 8-he...	25.18	9.1	ng	86147	6	24.79	189740	20.0
unknown25.92	25.92	6.2	ng	59177	6	24.79	189740	20.0
1,3a-Ethano-3aH-i...	26.02	7.7	ng	73259	6	24.79	189740	20.0
unknown26.20	26.20	10.7	ng	101568	6	24.79	189740	20.0
unknown26.47	26.47	4.7	ng	44629	6	24.79	189740	20.0
unknown26.88	26.88	4.7	ng	44910	6	24.79	189740	20.0
unknown27.54	27.54	6.0	ng	57357	6	24.79	189740	20.0
unknown27.80	27.80	14.6	ng	138313	6	24.79	189740	20.0