

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF092115\
 Data File : BF081736.D
 Acq On : 22 Sep 2015 6:16
 Operator : UM/IZ
 Sample : G3747-07
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TRENCH-2-3

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	885473	939659	41.89%	6.593%
2	1.281	15	17	29	rVB4	14861	37473	1.67%	0.263%
3	1.441	29	31	37	rBV	66234	160639	7.16%	1.127%
4	1.521	37	38	77	rVB	625421	2242894	100.00%	15.736%
5	4.333	281	284	289	rBV	13730	29225	1.30%	0.205%
6	4.424	289	292	312	rVB	201177	582456	25.97%	4.086%
7	5.316	367	370	396	rBV	495488	1003492	44.74%	7.040%
8	7.590	566	569	574	rBV	534770	997791	44.49%	7.000%
9	7.682	574	577	585	rVB	607214	1160472	51.74%	8.142%
10	8.162	615	619	625	rBB	130797	232217	10.35%	1.629%
11	8.482	644	647	657	rBB	495499	779639	34.76%	5.470%
12	9.465	729	733	745	rBV	383067	698790	31.16%	4.903%
13	11.054	869	872	879	rBV	164093	286352	12.77%	2.009%
14	13.705	1100	1104	1107	rBV	679896	1143435	50.98%	8.022%
15	13.762	1107	1109	1116	rVB	16385	37784	1.68%	0.265%
16	14.768	1192	1197	1204	rBV	79704	159881	7.13%	1.122%
17	15.191	1230	1234	1242	rVB	170637	310840	13.86%	2.181%
18	16.586	1352	1356	1360	rBV	19586	32462	1.45%	0.228%
19	17.100	1397	1401	1410	rBV	341016	706235	31.49%	4.955%
20	17.729	1452	1456	1462	rBV2	25879	79193	3.53%	0.556%
21	18.700	1537	1541	1545	rBV	170237	318597	14.20%	2.235%
22	18.814	1548	1551	1553	rVV	21595	36851	1.64%	0.259%
23	18.872	1553	1556	1566	rVB3	11225	27912	1.24%	0.196%
24	19.660	1619	1625	1629	rBV	28504	57689	2.57%	0.405%
25	19.843	1637	1641	1644	rBV	17072	29939	1.33%	0.210%
26	20.460	1692	1695	1701	rBV	28397	58735	2.62%	0.412%
27	21.409	1775	1778	1784	rVV2	31664	61654	2.75%	0.433%
28	21.740	1804	1807	1810	rBV	14418	25142	1.12%	0.176%
29	22.072	1832	1836	1838	rBV	1016942	1178028	52.52%	8.265%
30	23.318	1942	1945	1946	rBV	102826	164228	7.32%	1.152%
31	23.352	1946	1948	1955	rVV	208566	307338	13.70%	2.156%
32	24.003	2003	2005	2009	rVB2	21270	28517	1.27%	0.200%
33	24.358	2034	2036	2039	rVB	36021	53769	2.40%	0.377%
34	24.598	2055	2057	2061	rBV2	20370	33954	1.51%	0.238%

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ALS Vial : 24 Sample Multiplier: 1

Instrument :
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ClientSampleId :
TRENCH-2-3

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	24.792	2071	2074	2080	rVB	138402	215445	9.61%	1.512%
36	25.546	2136	2140	2143	rVB4	11625	34455	1.54%	0.242%

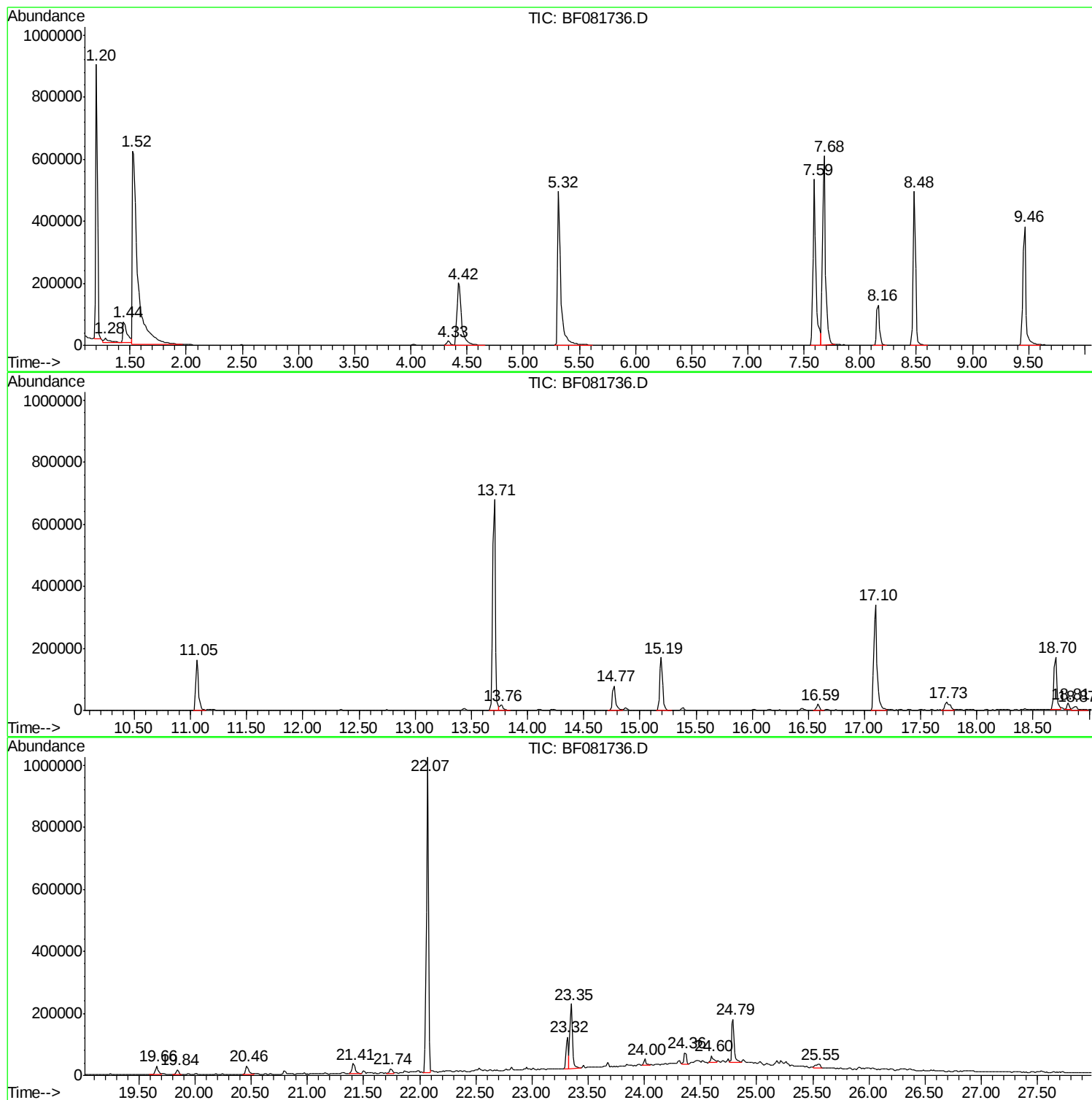
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF090115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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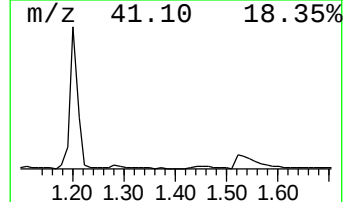
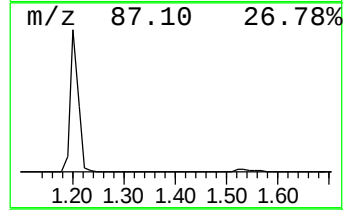
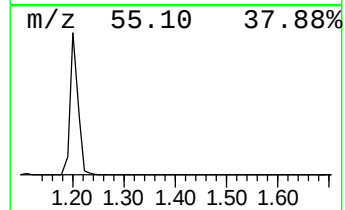
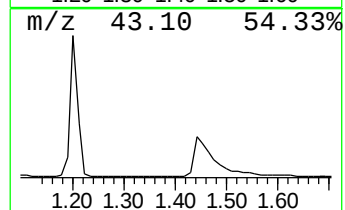
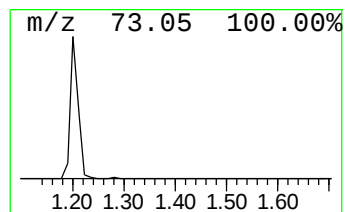
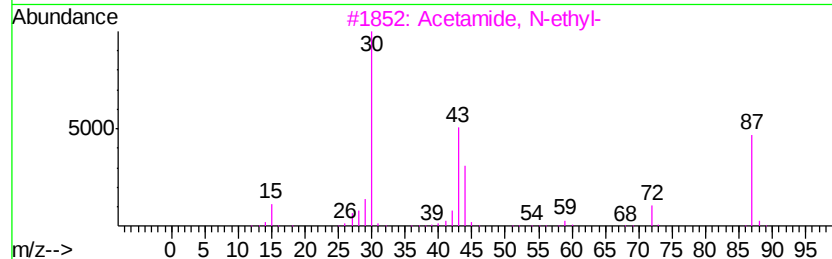
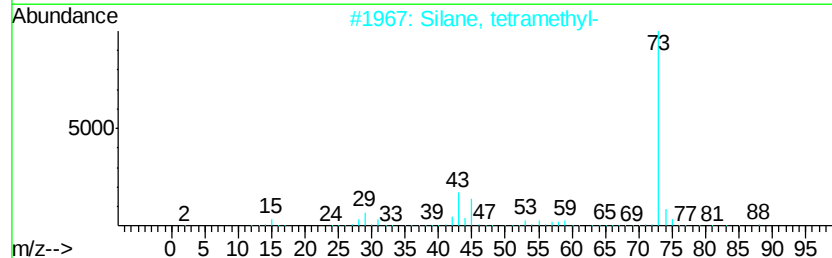
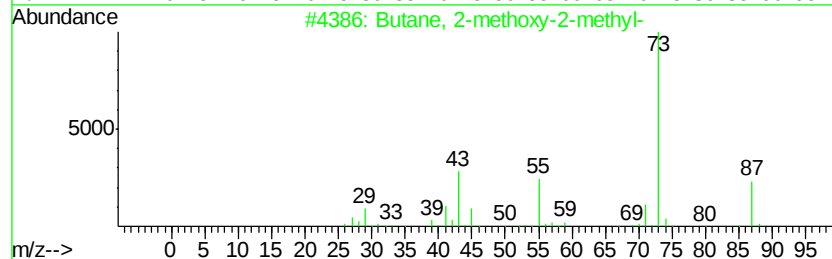
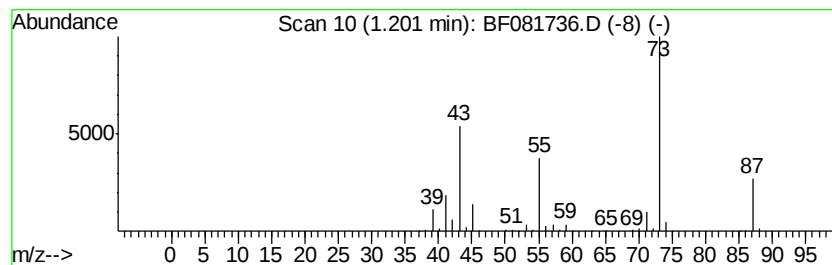
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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	80.93 ng	939659	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	Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9
4	1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9
5	1-Propene, 3-methoxy-	72	C4H8O	000627-40-7	9



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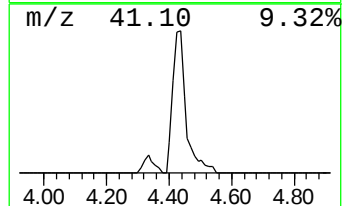
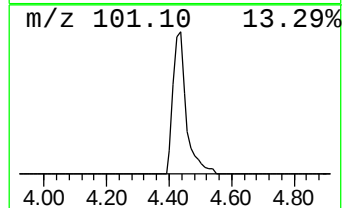
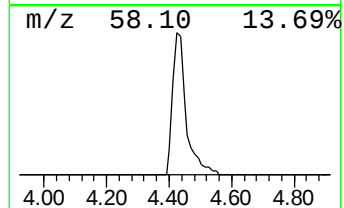
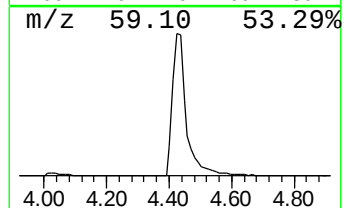
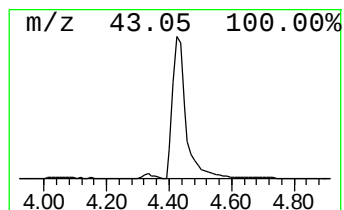
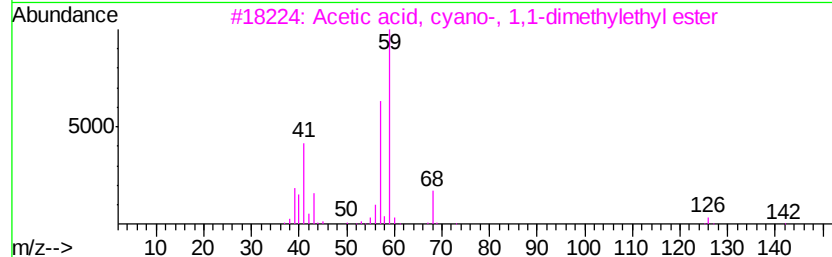
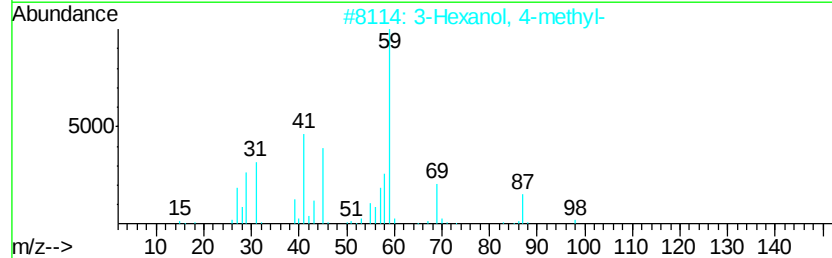
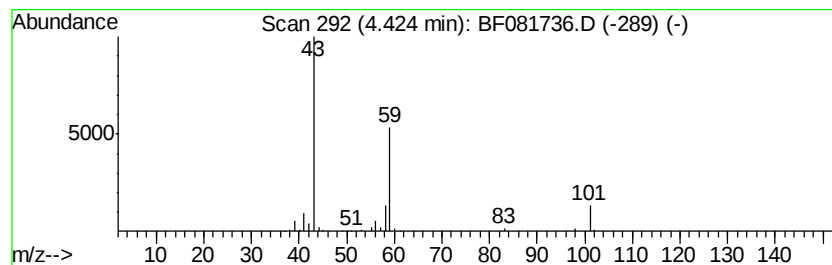
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF090115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.42	50.16 ng	582456	1,4-Dichlorobenzene-d4	8.16

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50	
2	3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33	
3	Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	25	
4	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10	
5	2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	9	



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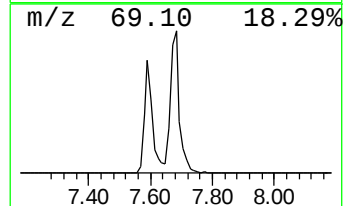
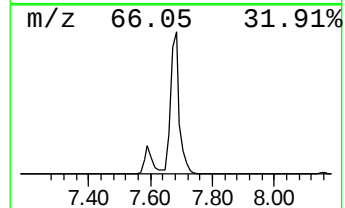
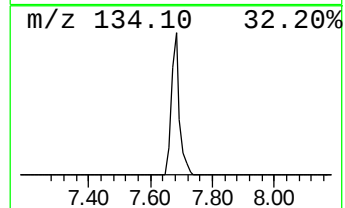
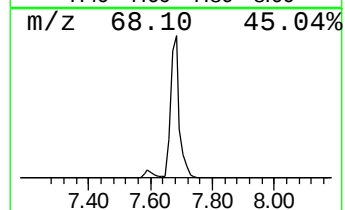
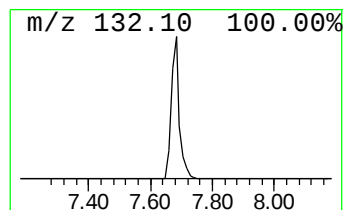
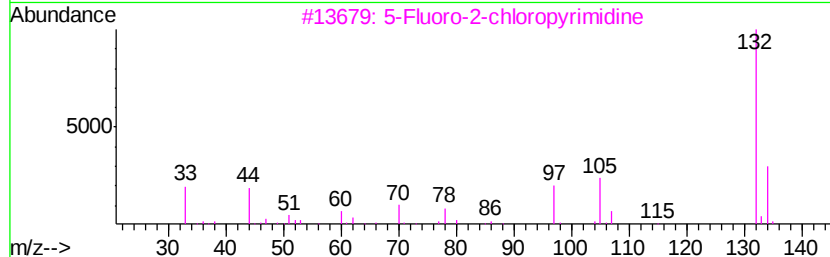
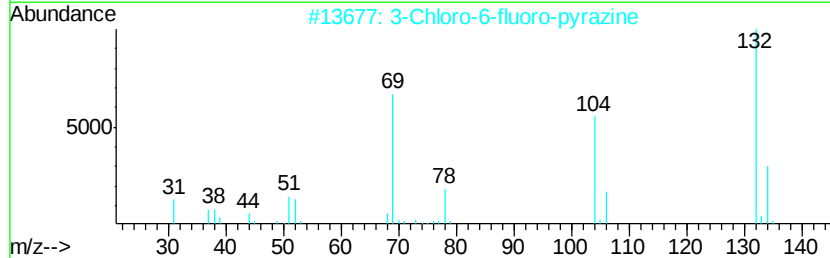
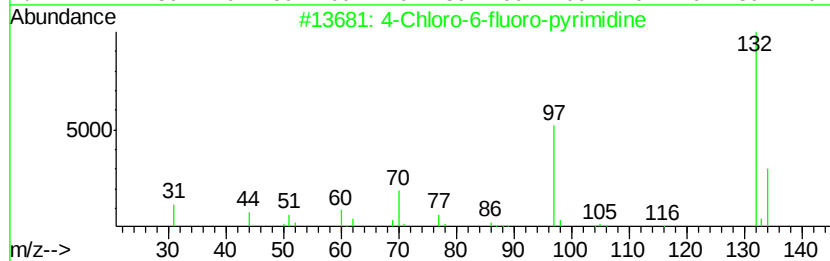
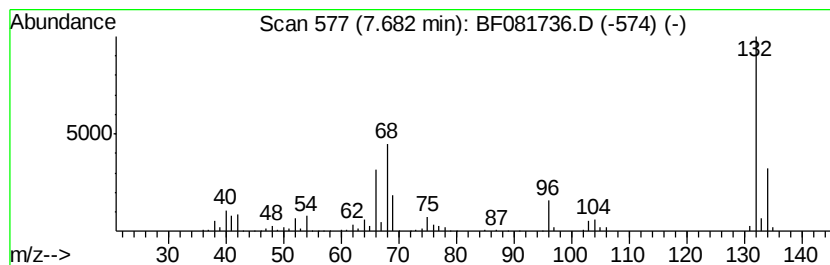
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 unknown7.68 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.68	99.95 ng	1160470	1,4-Dichlorobenzene-d4	8.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
3		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
4		(5-METHYL-2-PYRIDYL)ACETONITRILE	132	C8H8N2	1000241-93-9	10
5		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9



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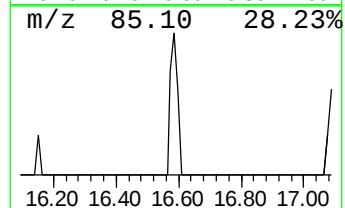
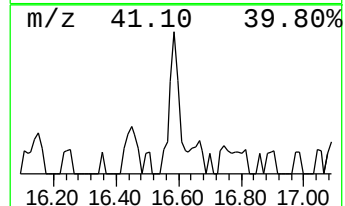
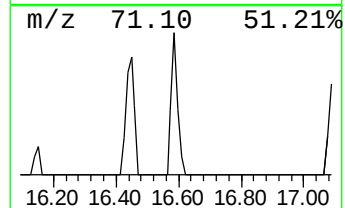
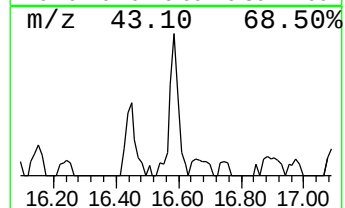
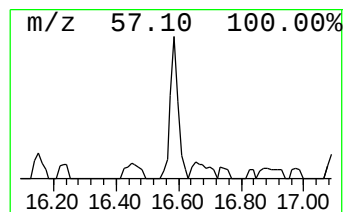
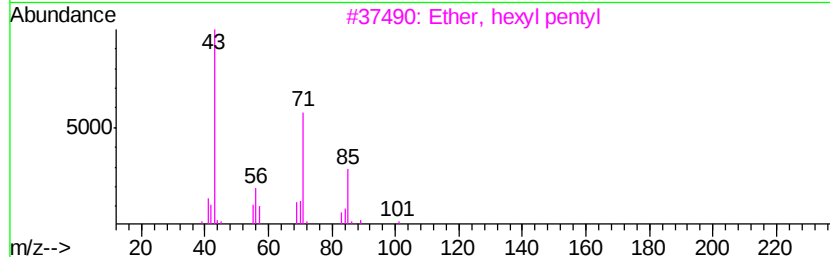
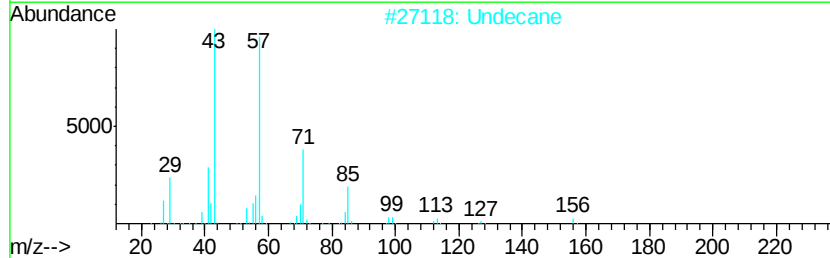
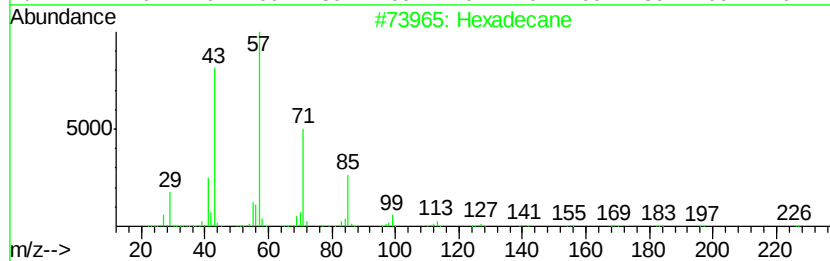
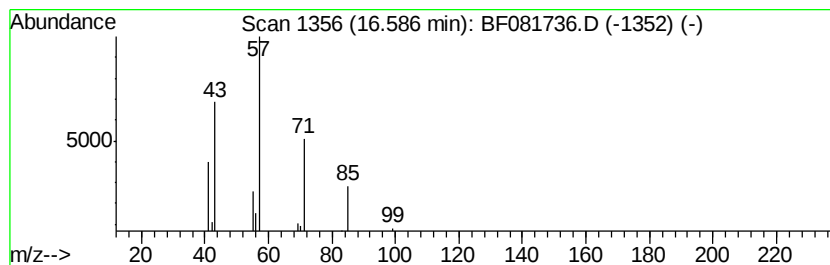
Instrument :
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ClientSampleId :
TRENCH-2-3

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 10 Hexadecane Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD		R.T.		
16.59	2.09 ng	32462	Acenaphthene-d10		15.19		
Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecane	226	C16H34	000544-76-3	83
2			Undecane	156	C11H24	001120-21-4	72
3			Ether, hexyl pentyl	172	C11H24O	032357-83-8	50
4			Tridecane	184	C13H28	000629-50-5	50
5			Hexane, 2,2,3,3-tetramethyl-	142	C10H22	013475-81-5	43



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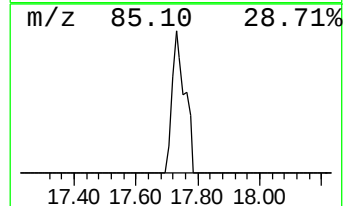
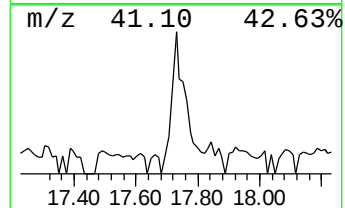
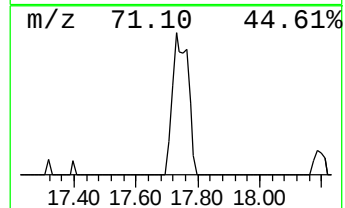
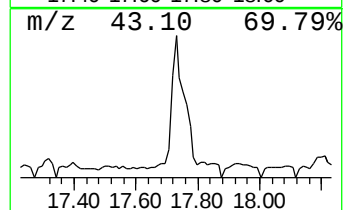
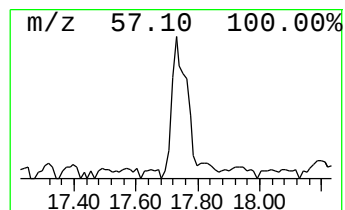
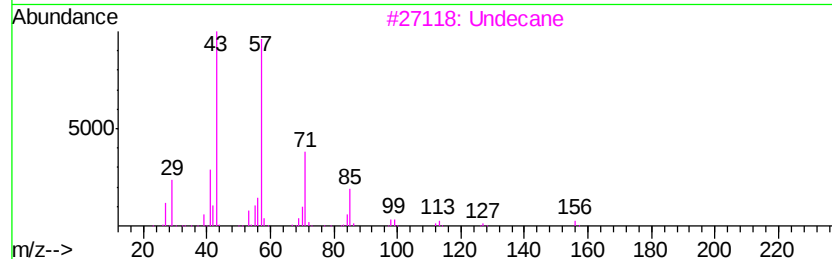
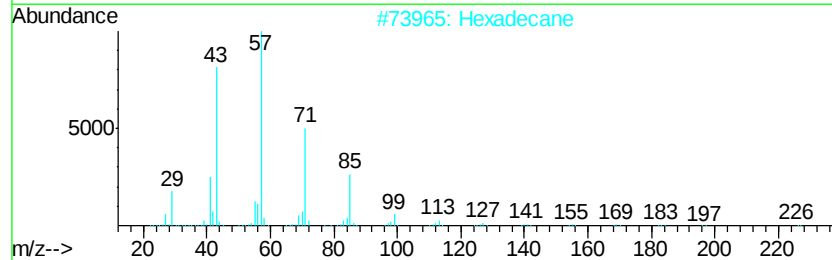
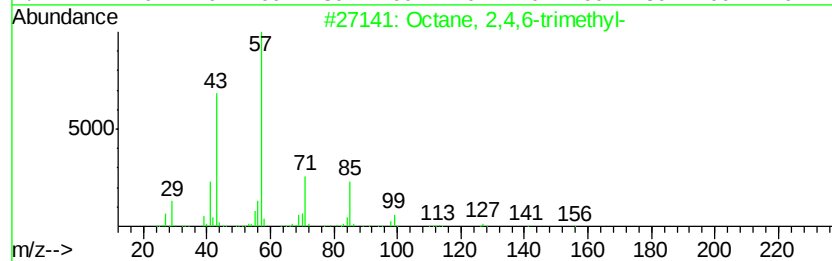
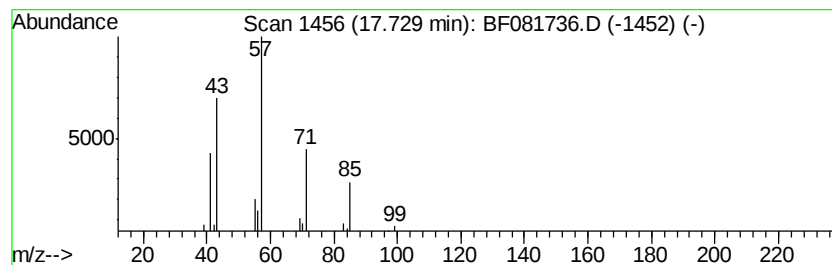
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 Octane, 2,4,6-trimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.73	4.97 ng	79193	Phenanthrene-d10	18.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octane, 2,4,6-trimethyl-	156	C11H24	062016-37-9	78
2		Hexadecane	226	C16H34	000544-76-3	78
3		Undecane	156	C11H24	001120-21-4	53
4		Tridecane	184	C13H28	000629-50-5	45
5		Ether, hexyl pentyl	172	C11H24O	032357-83-8	42



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ClientSampleId :
TRENCH-2-3

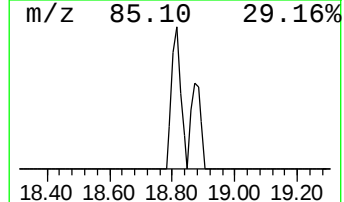
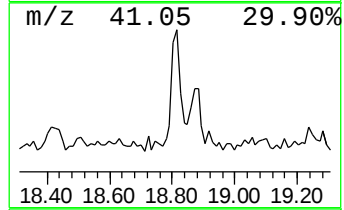
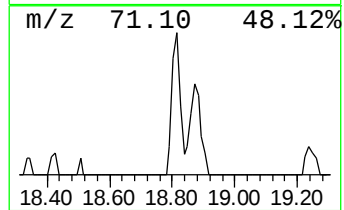
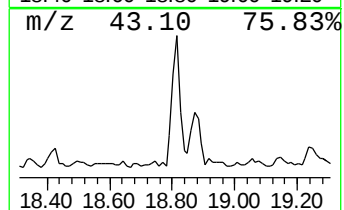
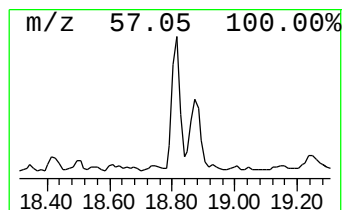
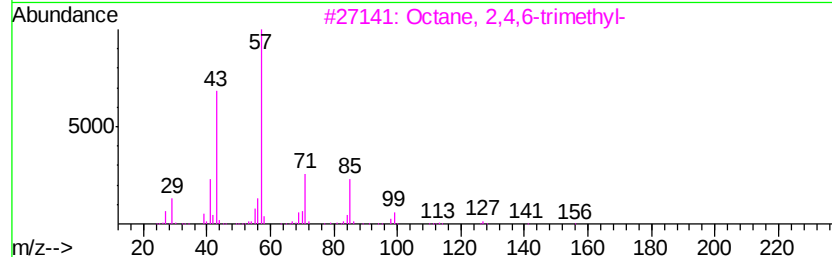
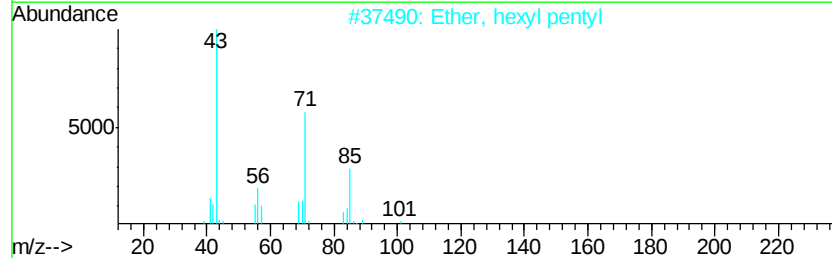
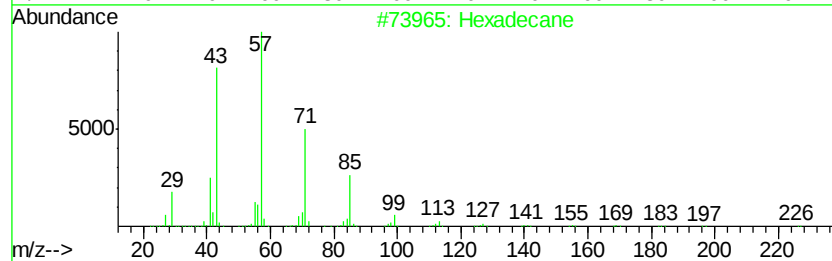
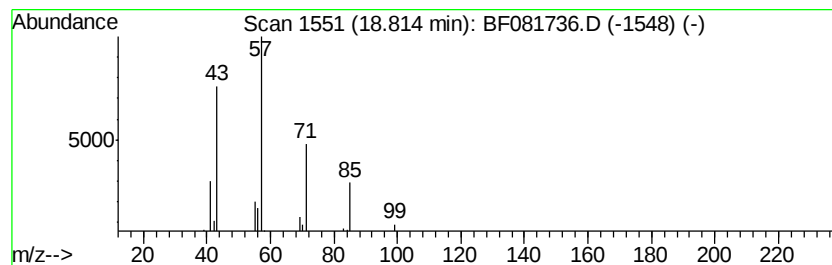
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Ether, hexyl pentyl Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.81	2.31 ng	36851	Phenanthrene-d10	18.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecane	226	C16H34	000544-76-3	83
2		Ether, hexyl pentyl	172	C11H24O	032357-83-8	72
3		Octane, 2,4,6-trimethyl-	156	C11H24	062016-37-9	64
4		Decane, 2,2,7-trimethyl-	184	C13H28	062237-99-4	64
5		Octane, 2-methyl-	128	C9H20	003221-61-2	59



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF092115\
Data File : BF081736.D
Acq On : 22 Sep 2015 6:16
Operator : UM/IZ
Sample : G3747-07
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TRENCH-2-3

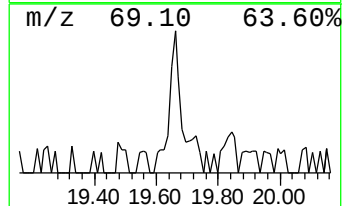
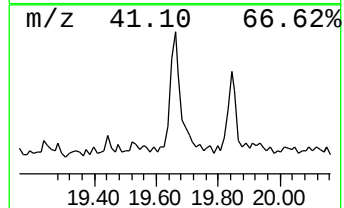
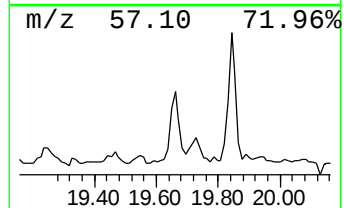
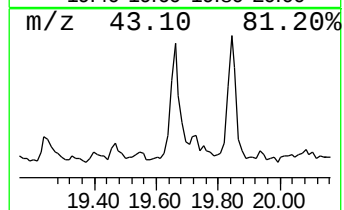
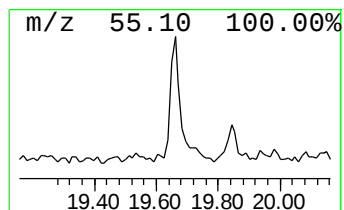
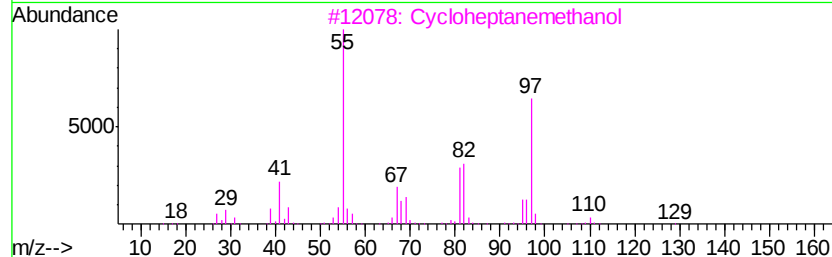
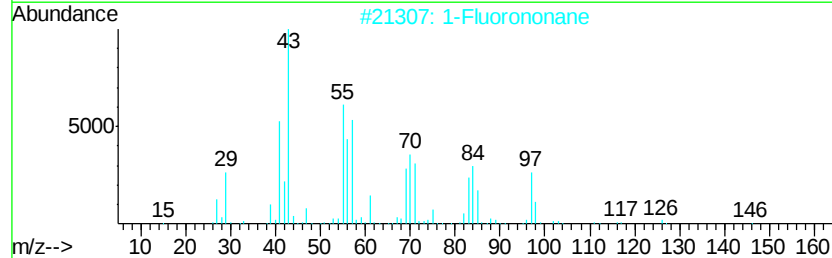
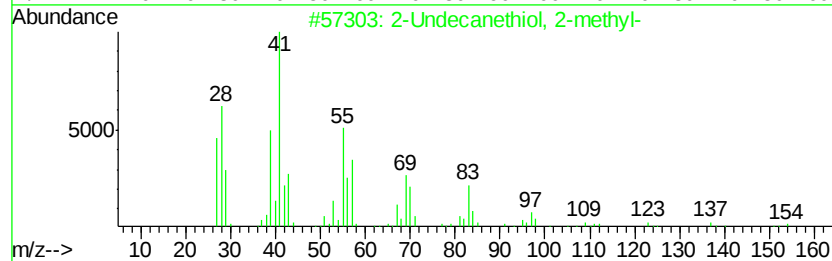
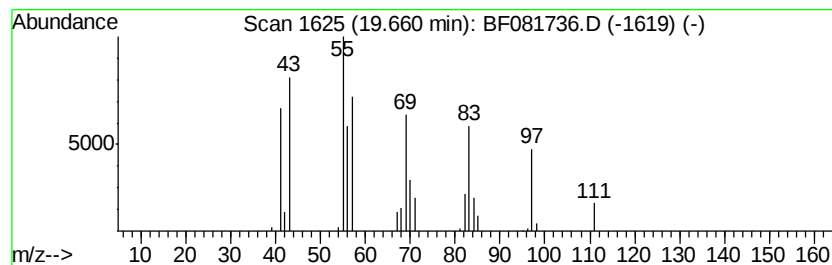
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 unknown19.66 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.66	3.62 ng	57689	Phenanthrene-d10	18.70

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Undecanethiol, 2-methyl-	202	C12H26S	010059-13-9	38		
2	1-Fluorononane	146	C9H19F	000463-18-3	35		
3	Cycloheptanemethanol	128	C8H16O	004448-75-3	27		
4	1-Heptanol, 6-methyl-	130	C8H18O	001653-40-3	27		
5	Isooctanol	130	C8H18O	026952-21-6	27		



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF092115\
Data File : BF081736.D
Acq On : 22 Sep 2015 6:16
Operator : UM/IZ
Sample : G3747-07
Misc :
ALS Vial : 24 Sample Multiplier: 1

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ClientSampleId :
TRENCH-2-3

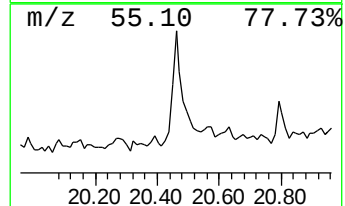
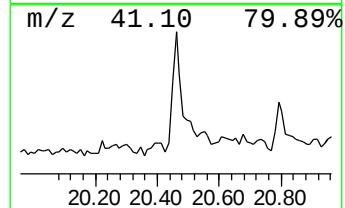
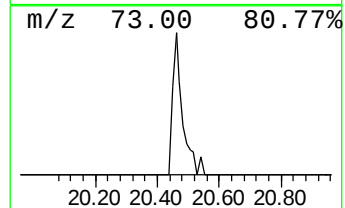
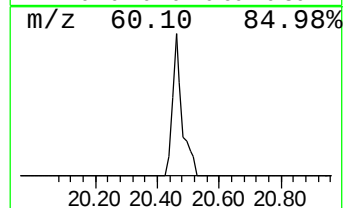
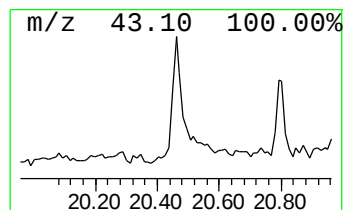
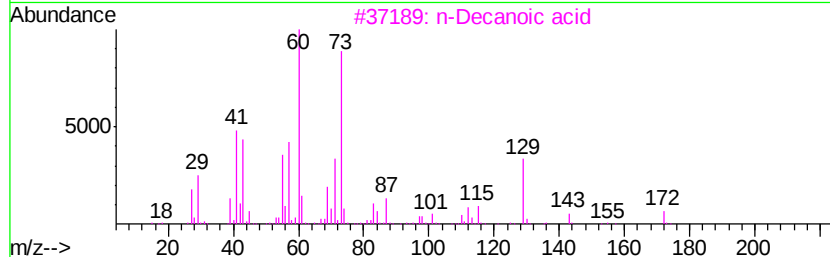
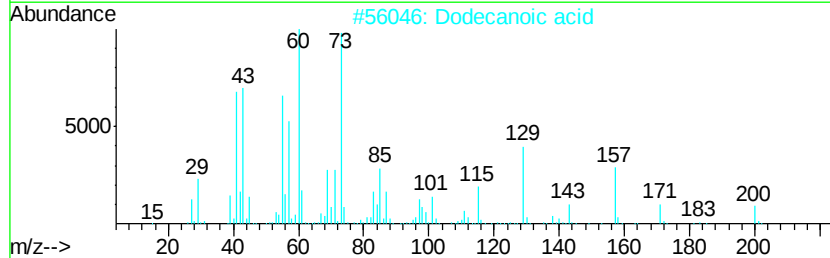
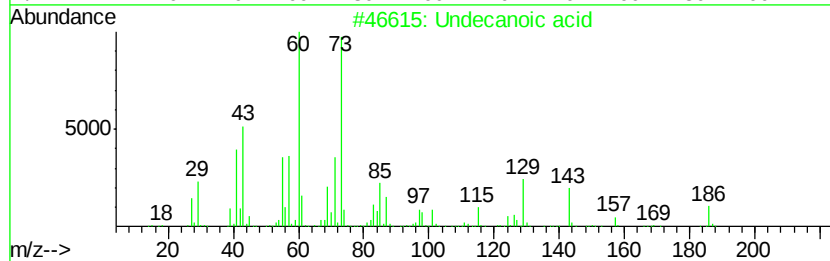
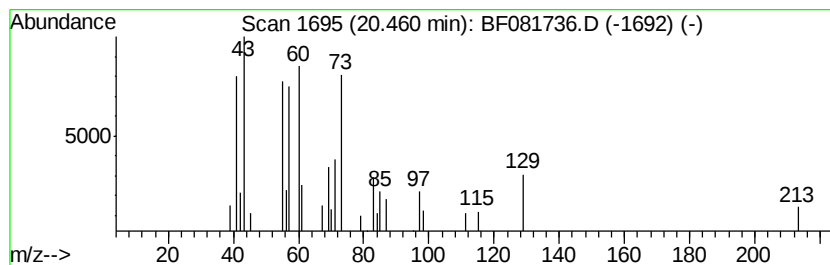
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Undecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.46	3.69 ng	58735	Phenanthrene-d10	18.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Undecanoic acid	186	C11H22O2	000112-37-8	64
2		Dodecanoic acid	200	C12H24O2	000143-07-7	59
3		n-Decanoic acid	172	C10H20O2	000334-48-5	50
4		4-Propionyloxvtridecane	256	C16H32O2	1000245-62-6	22
5		Amyl Nitrite	117	C5H11NO2	000110-46-3	22



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF092115\
Data File : BF081736.D
Acq On : 22 Sep 2015 6:16
Operator : UM/IZ
Sample : G3747-07
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TRENCH-2-3

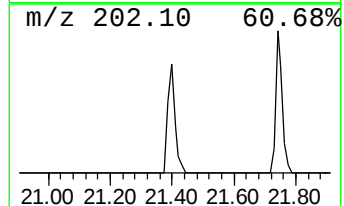
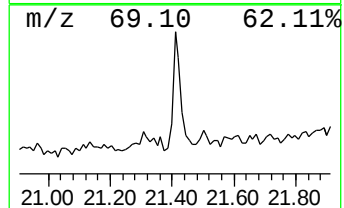
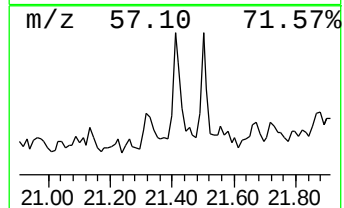
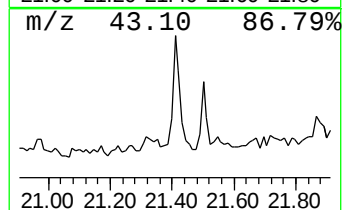
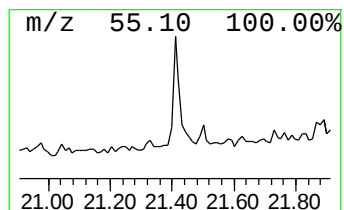
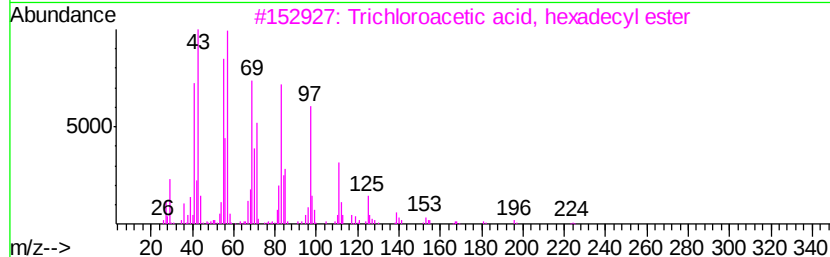
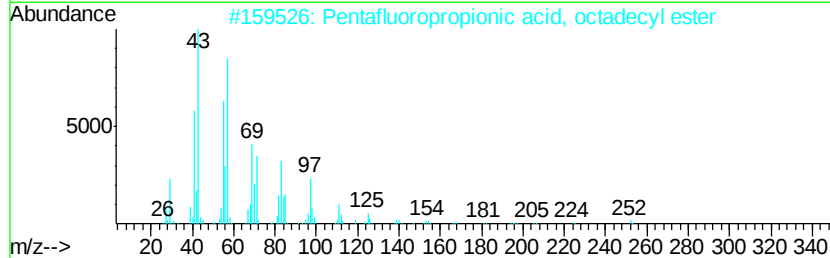
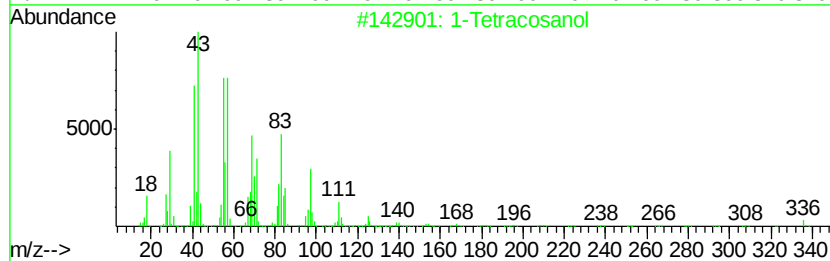
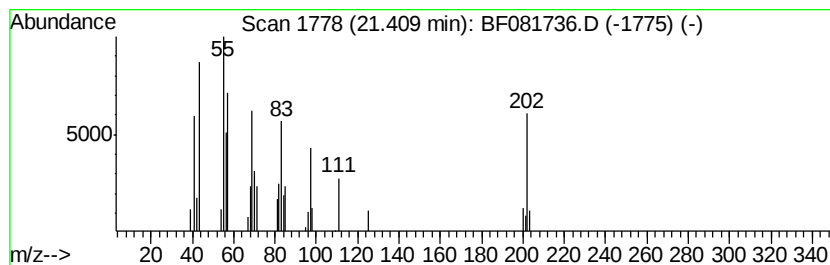
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 1-Tetracosanol Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.41	4.01 ng	61654	Chrysene-d12	23.35

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Tetracosanol	354	C24H50O	000506-51-4	62
2			Pentafluoropropionic acid, octadecyl ester	416	C21H37F5O2	1000280-07-7	58
3			Trichloroacetic acid, hexadecyl ester	386	C18H33Cl3O2	074339-54-1	58
4			3-Chloropropionic acid, heptadecyl ester	346	C20H39ClO2	1000283-05-1	53
5			1-Nonadecene	266	C19H38	018435-45-5	53



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF092115\
Data File : BF081736.D
Acq On : 22 Sep 2015 6:16
Operator : UM/IZ
Sample : G3747-07
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TRENCH-2-3

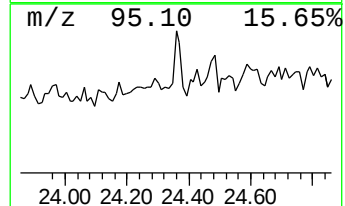
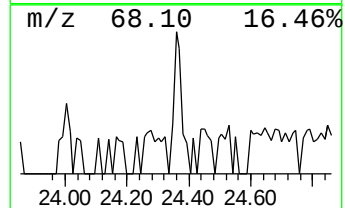
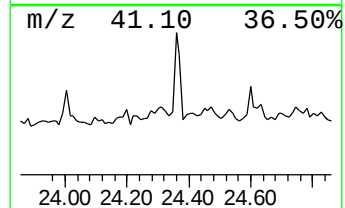
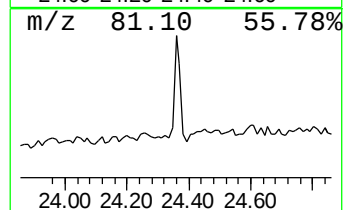
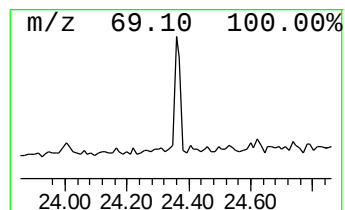
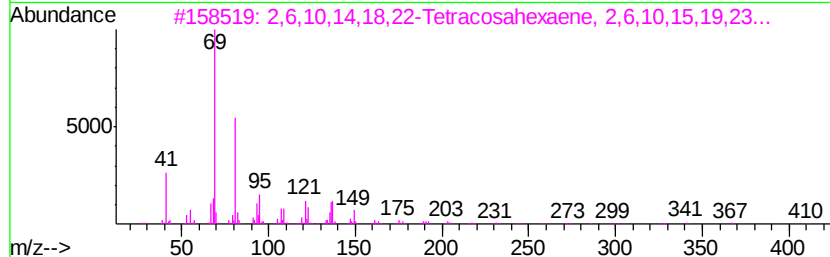
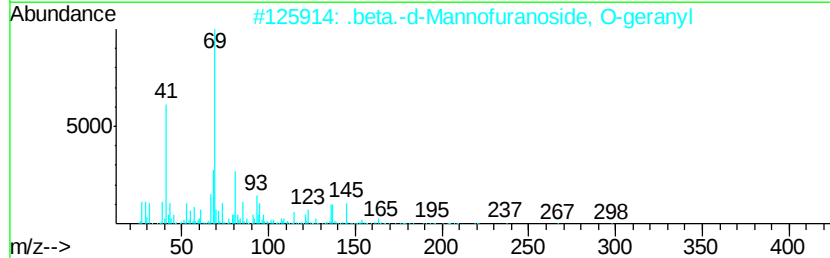
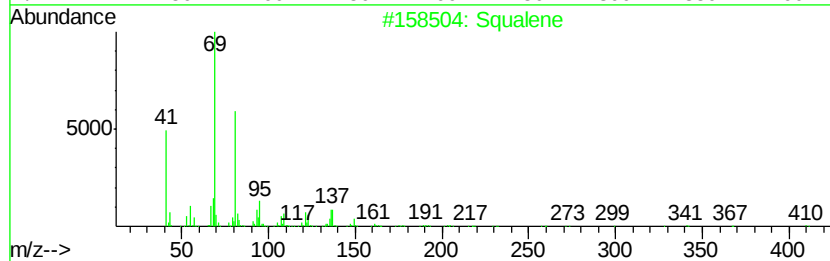
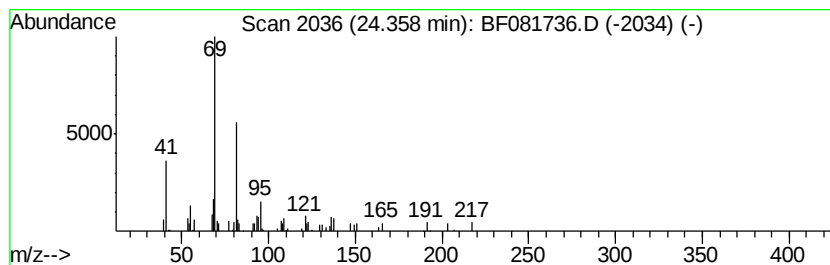
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Squalene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.36	4.99 ng	53769	Perylene-d12	24.79

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Squalene	410	C30H50	007683-64-9	64
2		.beta.-d-Mannofuranoside, O-geranyl	316	C16H28O6	1000154-77-2	59
3		2,6,10,14,18,22-Tetracosahexaene...	410	C30H50	000111-02-4	59
4		7,11-Dimethyldodeca-2,6,10-trien...	208	C14H24O	125529-10-4	50
5		2,6,10-Dodecatrien-1-ol, 3,7,11-...	222	C15H26O	004602-84-0	45



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF092115\
Data File : BF081736.D
Acq On : 22 Sep 2015 6:16
Operator : UM/IZ
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Misc :
ALS Vial : 24 Sample Multiplier: 1

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ClientSampleId :
TRENCH-2-3

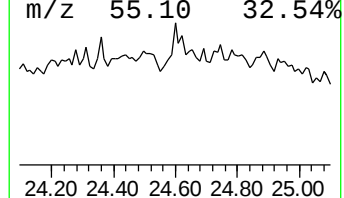
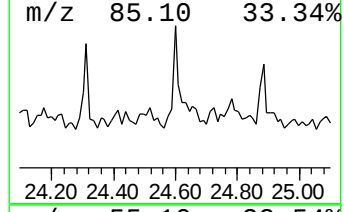
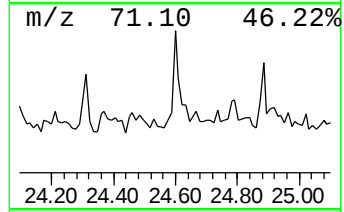
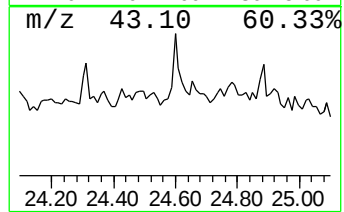
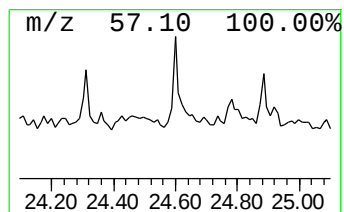
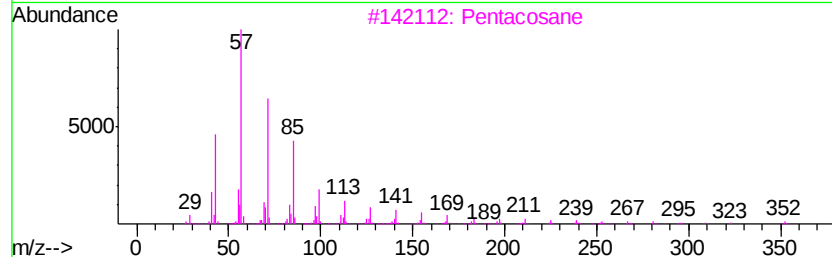
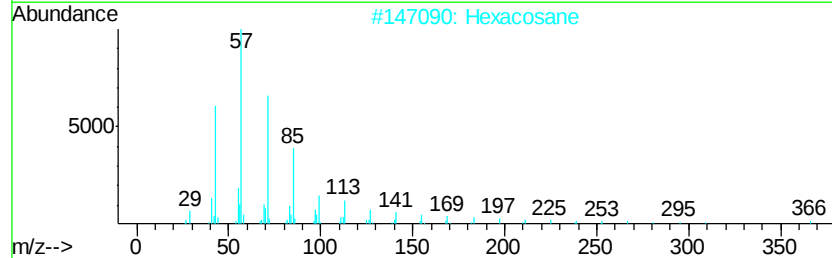
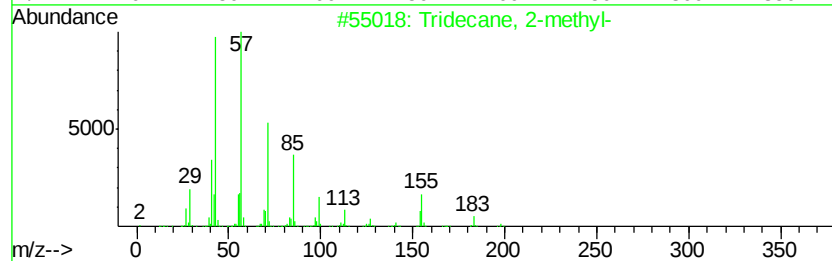
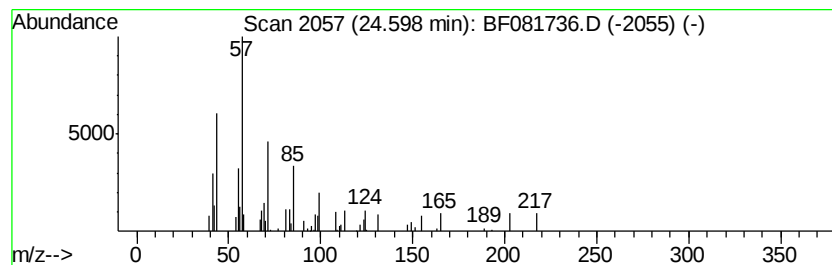
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 Tridecane, 2-methyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.60	3.15 ng	33954	Perylene-d12	24.79

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecane, 2-methyl-	198	C14H30	001560-96-9	53
2			Hexacosane	366	C26H54	000630-01-3	52
3			Pentacosane	352	C25H52	000629-99-2	52
4			Octadecane, 1-iodo-	380	C18H37I	000629-93-6	50
5			Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	50



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF092115\
Data File : BF081736.D
Acq On : 22 Sep 2015 6:16
Operator : UM/IZ
Sample : G3747-07
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
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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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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2-Pentanone, 4-hy...	4.42	50.2	ng	582456	1	8.16	232217	20.0
unknown7.68	7.68	100.0	ng	1160470	1	8.16	232217	20.0
Hexadecane	16.59	2.1	ng	32462	3	15.19	310840	20.0
Octane, 2,4,6-tri...	17.73	5.0	ng	79193	4	18.70	318597	20.0
Ether, hexyl pentyl	18.81	2.3	ng	36851	4	18.70	318597	20.0
unknown19.66	19.66	3.6	ng	57689	4	18.70	318597	20.0
Undecanoic acid	20.46	3.7	ng	58735	4	18.70	318597	20.0
1-Tetracosanol	21.41	4.0	ng	61654	5	23.35	307338	20.0
Squalene	24.36	5.0	ng	53769	6	24.79	215445	20.0
Tridecane, 2-methyl-	24.60	3.1	ng	33954	6	24.79	215445	20.0