

Data Path : Z:\HPCHEM1\BNA F\DATA\BF051215\
 Data File : BF079174.D
 Acq On : 12 May 2015 23:06
 Operator : TP/IZ
 Sample : G2221-01
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 ALMASI-SOIL

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	1.438	19	22	24	rVB	3788442	4778731	100.00%	19.156%
2	1.564	30	33	38	rVB	71171	149173	3.12%	0.598%
3	1.724	45	47	56	rVV2	224545	539255	11.28%	2.162%
4	1.850	56	58	90	rVB	1531012	2749619	57.54%	11.022%
5	5.073	338	340	346	rVB	126310	154124	3.23%	0.618%
6	5.587	382	385	387	rBV	1254659	1299624	27.20%	5.210%
7	6.845	493	495	498	rBV	1203912	1279783	26.78%	5.130%
8	6.993	506	508	511	rBV	1398991	1370412	28.68%	5.494%
9	7.279	531	533	536	rBV	320190	314375	6.58%	1.260%
10	7.473	547	550	552	rBV	884203	1069259	22.38%	4.286%
11	7.976	591	594	596	rBV	867223	855709	17.91%	3.430%
12	8.856	669	671	674	rBV	502385	439595	9.20%	1.762%
13	10.205	786	789	791	rBV	1784794	1630339	34.12%	6.535%
14	10.708	831	833	836	rVB	70987	62900	1.32%	0.252%
15	11.028	859	861	864	rVB	393140	479817	10.04%	1.923%
16	12.011	944	947	950	rBV	1033760	1083551	22.67%	4.344%
17	12.879	1020	1023	1024	rBV	490917	509907	10.67%	2.044%
18	12.902	1024	1025	1028	rVB	58373	56256	1.18%	0.226%
19	14.388	1152	1155	1156	rBV	95461	112219	2.35%	0.450%
20	14.662	1177	1179	1182	rVB	125287	131790	2.76%	0.528%
21	14.868	1195	1197	1200	rBV	1314666	1682592	35.21%	6.745%
22	15.634	1256	1264	1267	rBV9	17078	82848	1.73%	0.332%
23	16.045	1297	1300	1305	rBV	313114	467020	9.77%	1.872%
24	16.171	1308	1311	1313	rVV	494919	705429	14.76%	2.828%
25	16.217	1313	1315	1320	rVB2	202201	278234	5.82%	1.115%
26	16.343	1324	1326	1332	rVV6	17517	52412	1.10%	0.210%
27	16.880	1371	1373	1378	rBV	200380	377707	7.90%	1.514%
28	17.166	1396	1398	1403	rBV5	24021	49683	1.04%	0.199%
29	17.451	1421	1423	1425	rBV	63603	87042	1.82%	0.349%
30	17.497	1425	1427	1433	rVB	61472	127849	2.68%	0.513%
31	17.680	1441	1443	1445	rBV	114975	213260	4.46%	0.855%
32	17.771	1449	1451	1454	rVV2	48754	76321	1.60%	0.306%
33	17.829	1454	1456	1458	rVV	48792	73315	1.53%	0.294%
34	17.886	1459	1461	1465	rVV	47240	91203	1.91%	0.366%

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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
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35	17.954	1465	1467	1474	rVB	455106	685416	14.34%	2.748%
36	18.286	1494	1496	1499	rBV	67709	111478	2.33%	0.447%
37	18.537	1515	1518	1520	rVB	104400	156251	3.27%	0.626%
38	18.583	1520	1522	1524	rBV	78159	132003	2.76%	0.529%
39	18.777	1537	1539	1544	rVB2	47086	84359	1.77%	0.338%
40	19.337	1585	1588	1592	rBV	90235	189086	3.96%	0.758%
41	20.046	1647	1650	1653	rBV3	62006	155984	3.26%	0.625%

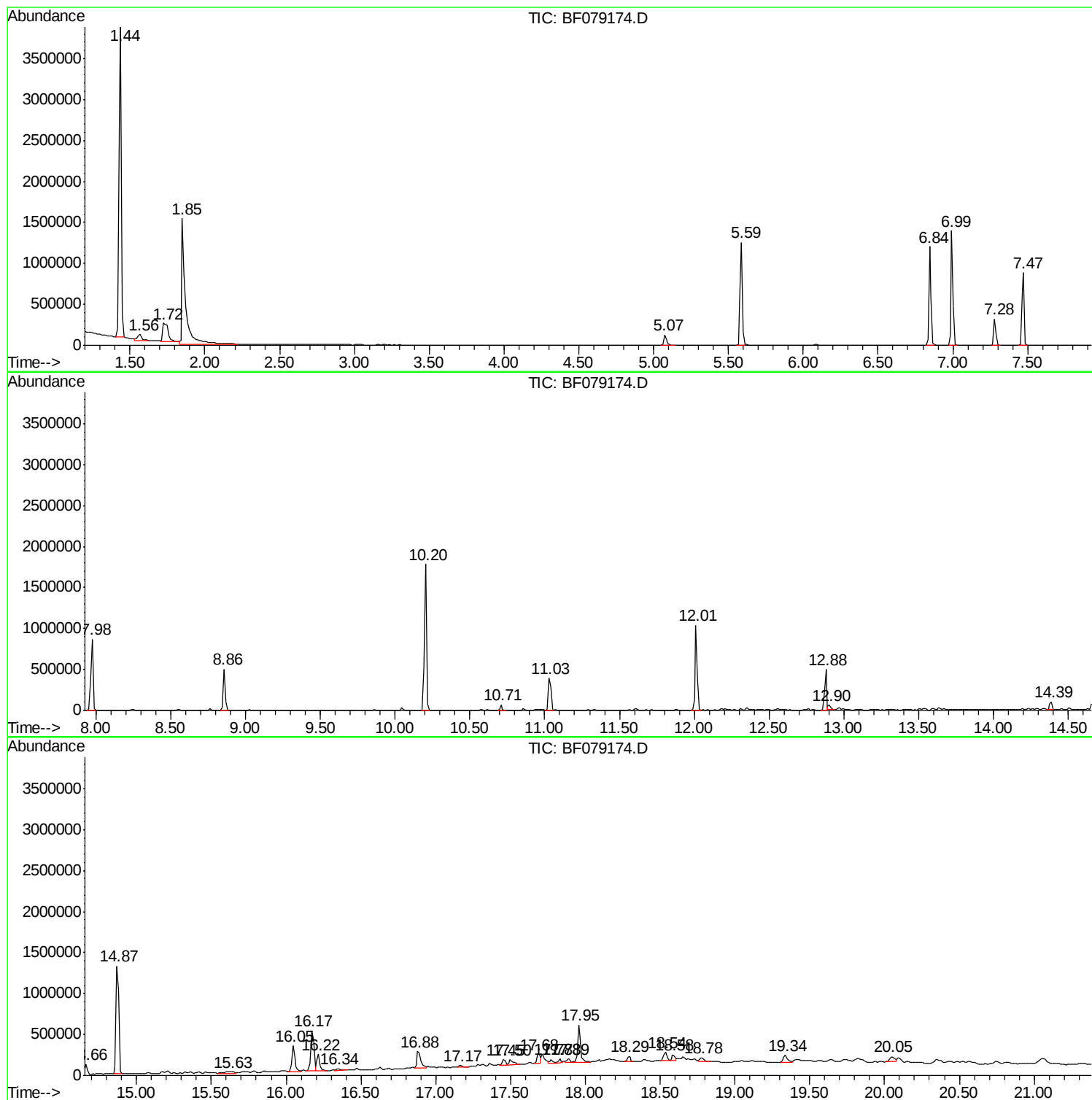
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF043015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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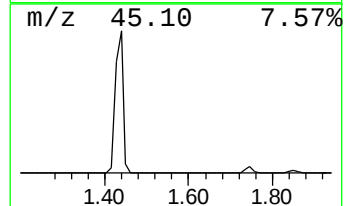
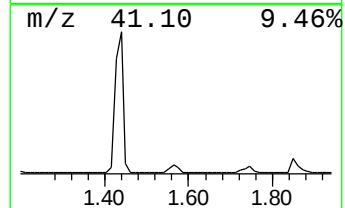
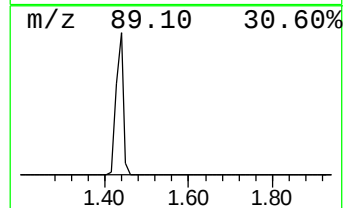
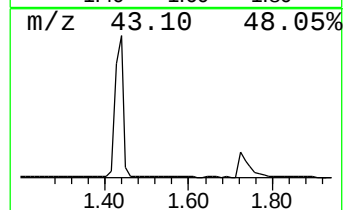
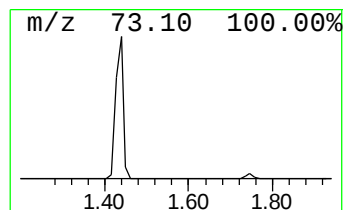
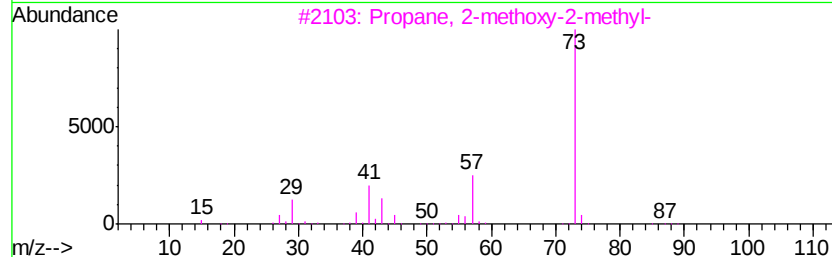
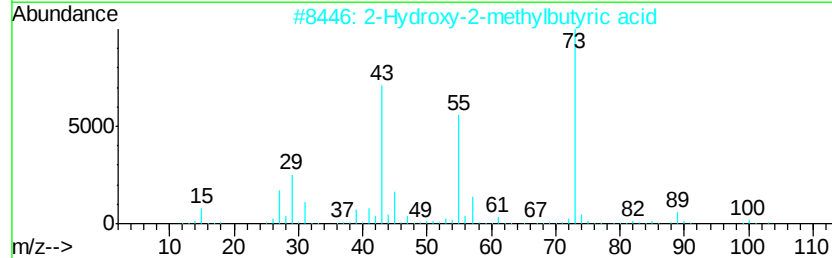
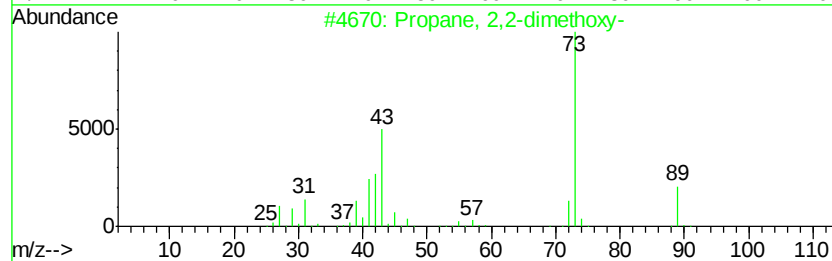
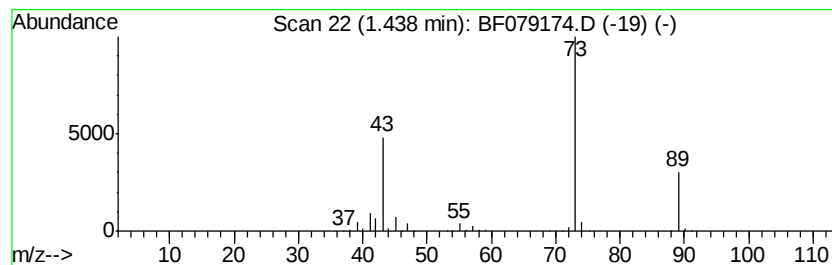
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown1.44 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.44	304.02 ng	4778730	1,4-Dichlorobenzene-d4	7.28

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propane, 2,2-dimethoxy-	104	C5H12O2	000077-76-9	38
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		1-Hexen-4-ol, 1-chloro-3,5-dimet...	162	C8H15ClO	100244-09-5	9
5		N-Ethylformamide	73	C3H7NO	000627-45-2	5



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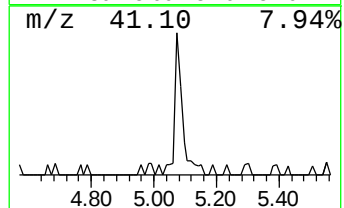
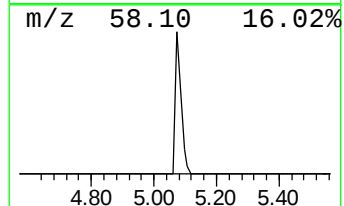
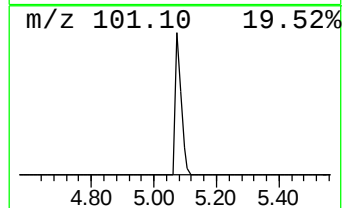
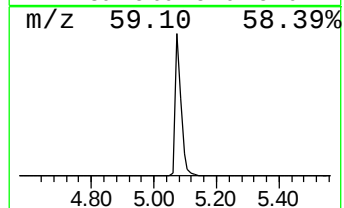
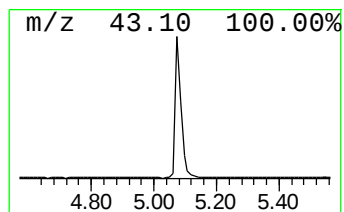
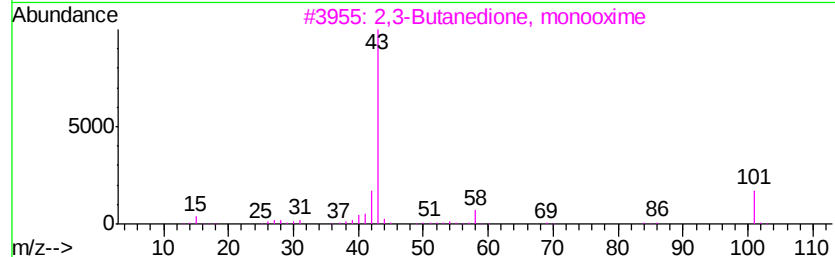
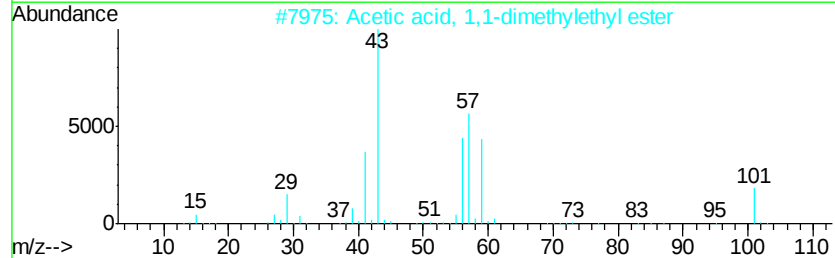
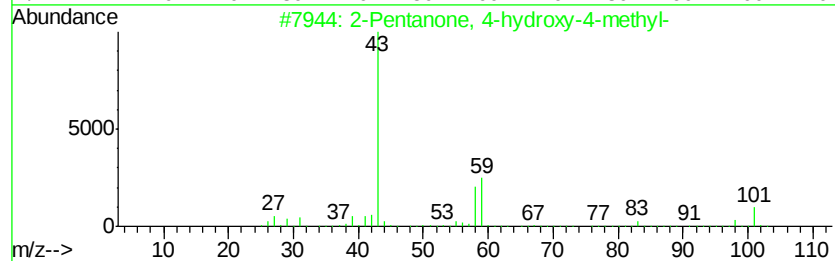
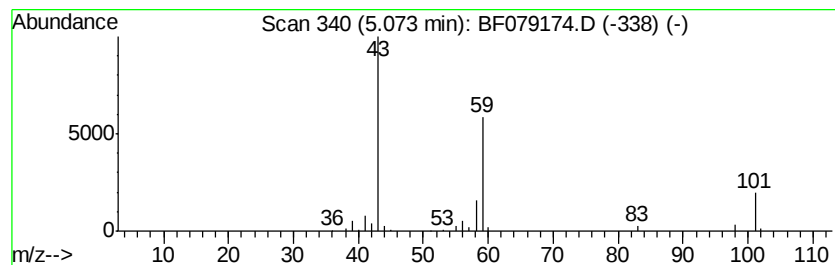
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 5 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.07	9.81 ng	154124	1,4-Dichlorobenzene-d4	7.28

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25
4		Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23
5		Guanidine	59	CH5N3	000113-00-8	9



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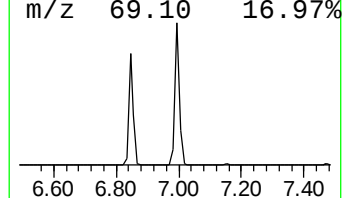
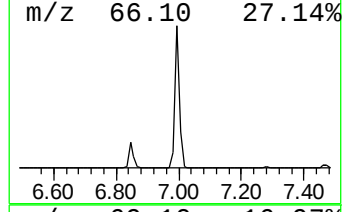
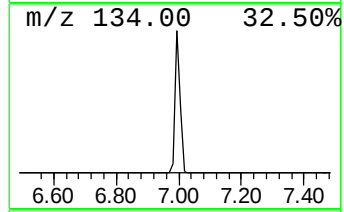
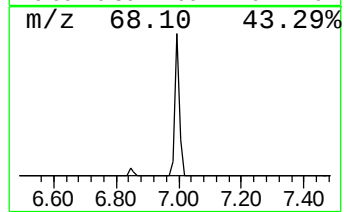
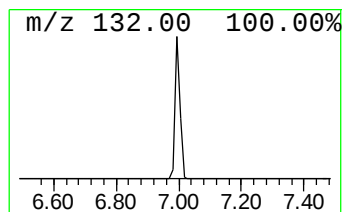
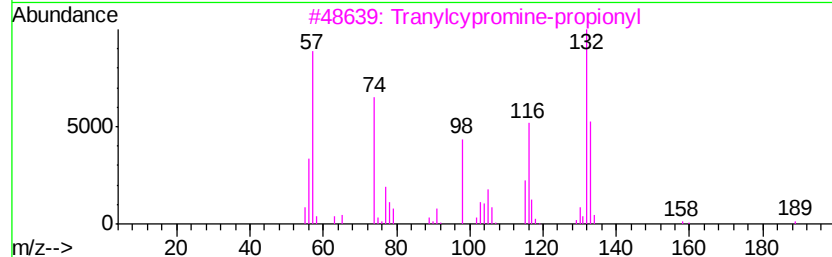
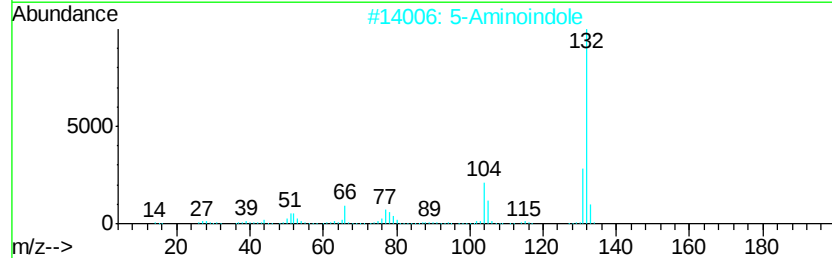
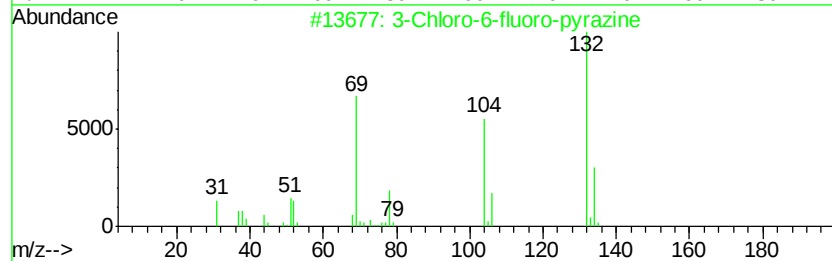
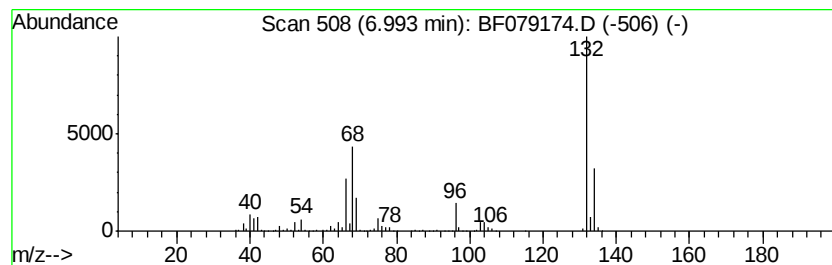
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 unknown6.99 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.99	87.18 ng	1370410	1,4-Dichlorobenzene-d4	7.28

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Chloro-6-fluoro-pyrazine			132	C4H2ClFN2	1000146-10-7	27
2	5-Aminoindole			132	C8H8N2	005192-03-0	12
3	Tranvlcvpromine-propionyl			189	C12H15NO	1000123-86-3	10
4	Bicyclo[4.2.0]octa-1,3,5-triene,...			132	C10H12	028749-81-7	9
5	4-Chloro-2-fluoro-pyrimidine			132	C4H2ClFN2	051422-00-5	9



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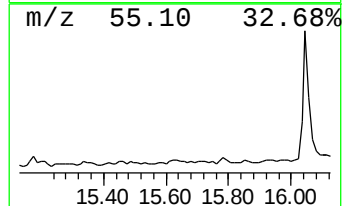
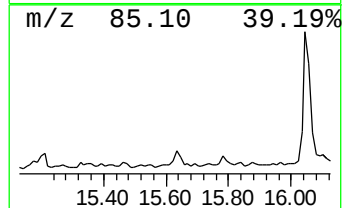
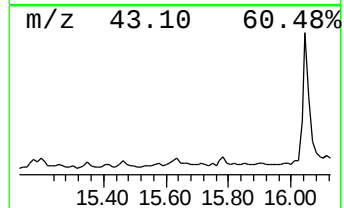
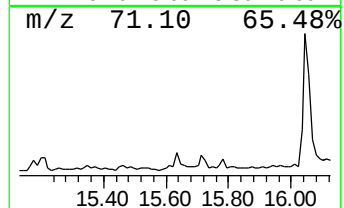
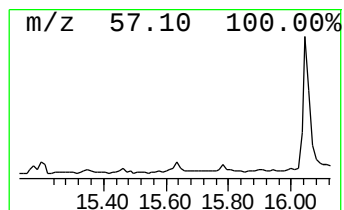
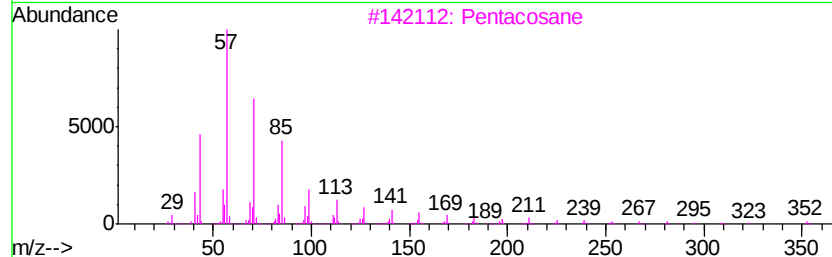
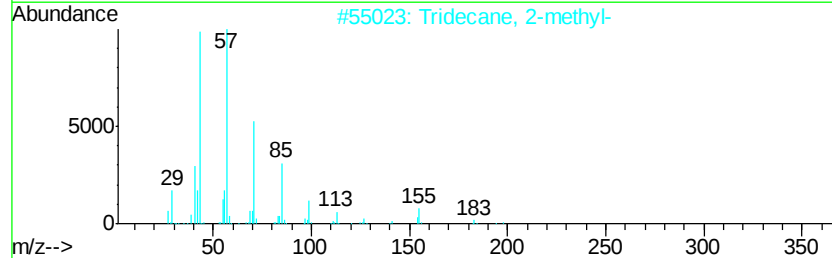
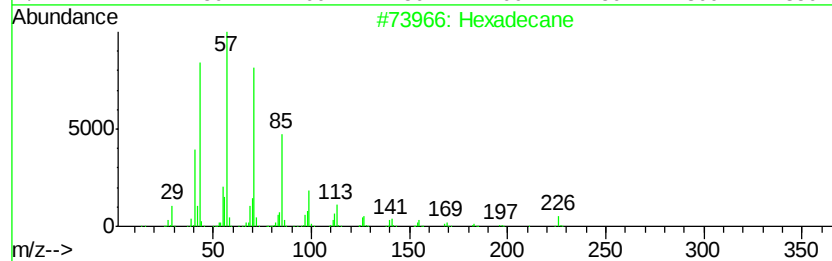
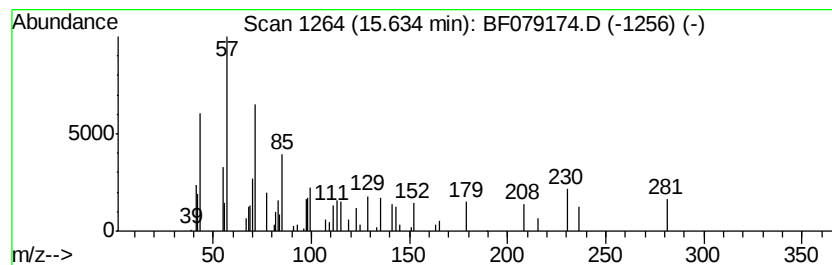
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 8 unknown15.63 Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.63	2.35 ng	82848	Chrysene-d12	16.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecane	226	C16H34	000544-76-3	43
2		Tridecane, 2-methyl-	198	C14H30	001560-96-9	43
3		Pentacosane	352	C25H52	000629-99-2	43
4		Pentadecane, 2,6,10-trimethyl-	254	C18H38	003892-00-0	43
5		Heptacosane, 1-chloro-	414	C27H55Cl	062016-79-9	38



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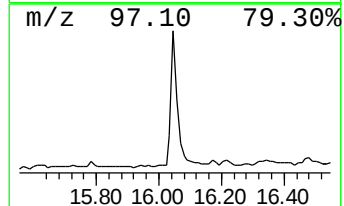
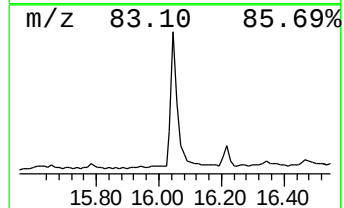
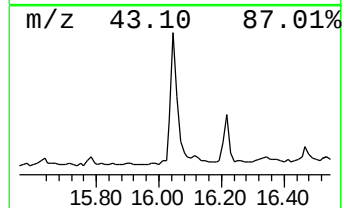
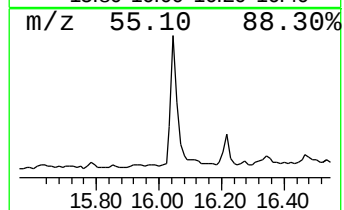
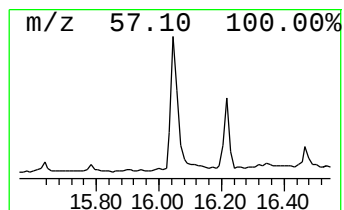
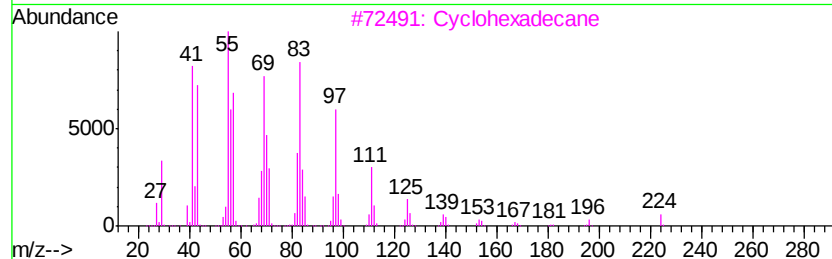
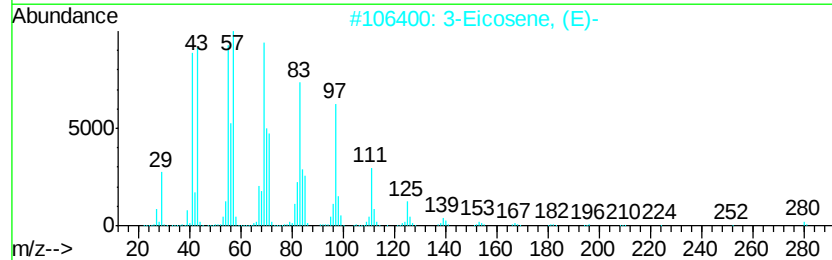
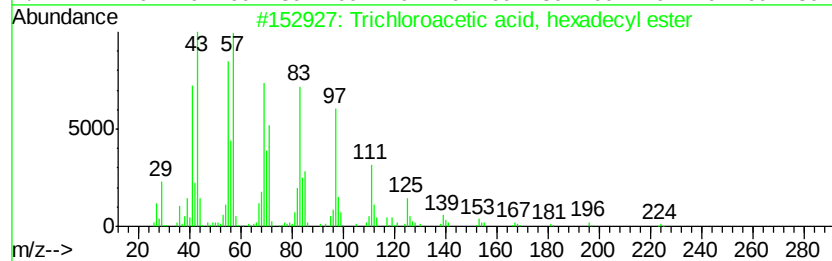
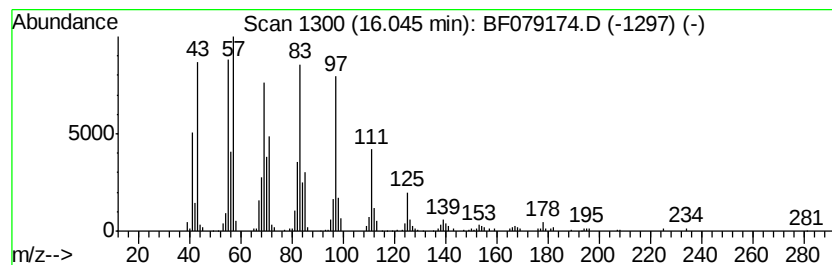
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Trichloroacetic acid, hexad... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.05	13.24 ng	467020	Chrysene-d12	16.17

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	91
2		3-Eicosene, (E)-	280	C20H40	074685-33-9	91
3		Cyclohexadecane	224	C16H32	000295-65-8	87
4		1-Heptadecanol	256	C17H36O	001454-85-9	87
5		Cyclopentadecane	210	C15H30	000295-48-7	87



Data Path : Z:\HPCHEM1\BNA F\DATA\BF051215\
Data File : BF079174.D
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Operator : TP/IZ
Sample : G2221-01
Misc :
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ClientSampleId :
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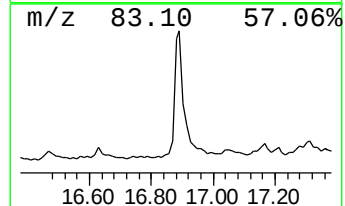
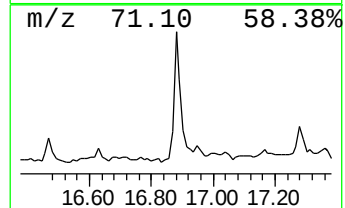
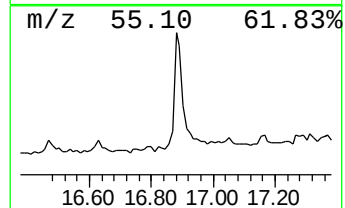
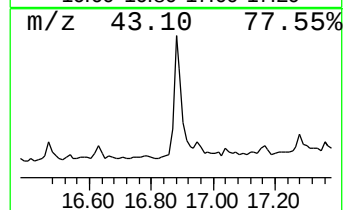
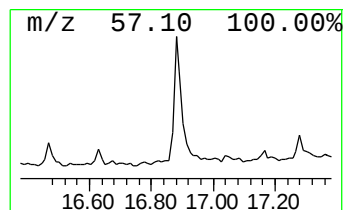
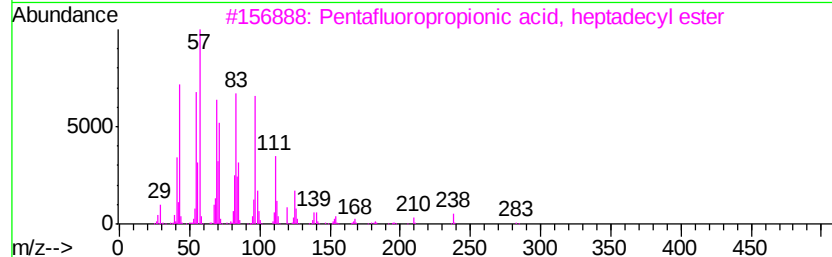
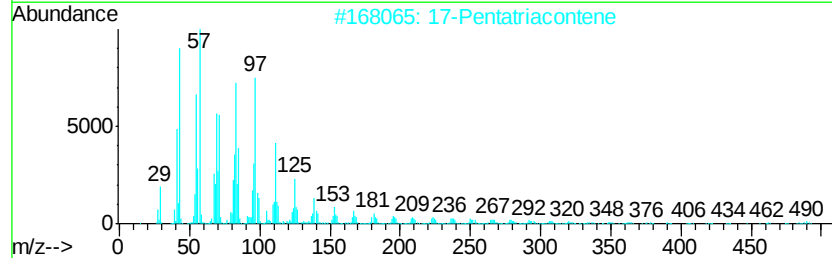
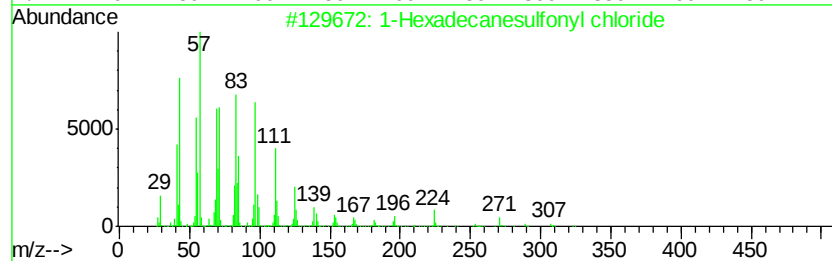
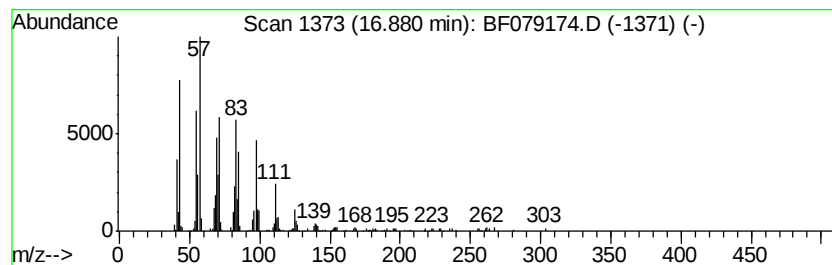
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 1-Hexadecanesulfonyl chloride Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.88	10.71 ng	377707	Chrysene-d12	16.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Hexadecanesulfonyl chloride	324	C16H33ClO2S	038775-38-1	87
2		17-Pentatriacontene	491	C35H70	006971-40-0	83
3		Pentafluoropropionic acid, hepta...	402	C20H35F5O2	1000283-04-2	76
4		1-Octadecanol	270	C18H38O	000112-92-5	70
5		Heptafluorobutanoic acid, heptad...	452	C21H35F7O2	1000282-97-3	70



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051215\
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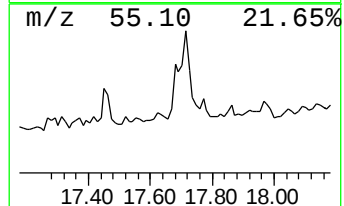
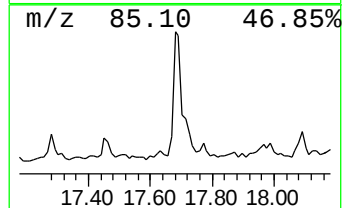
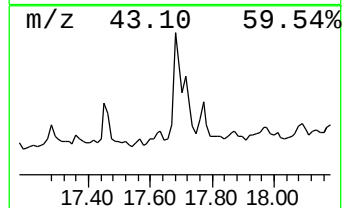
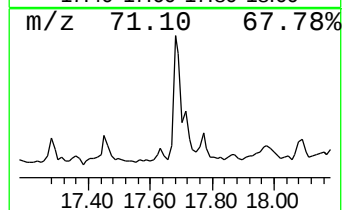
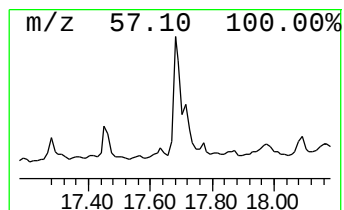
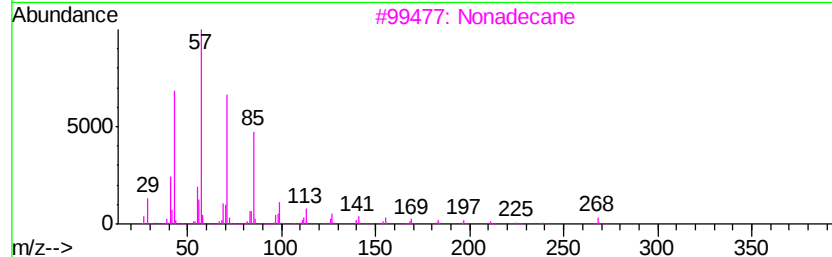
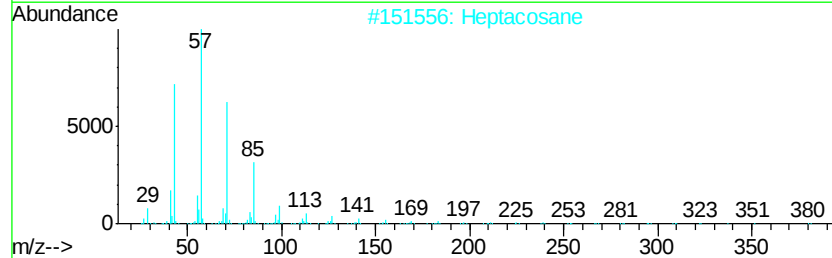
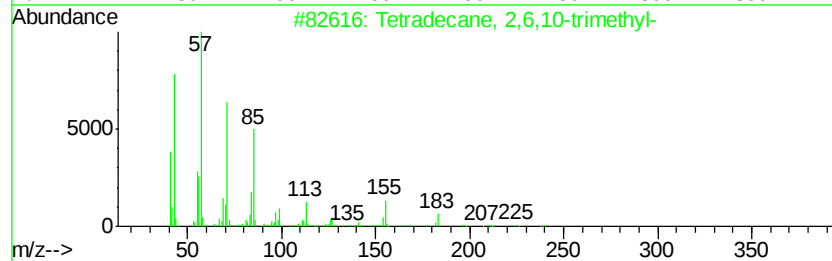
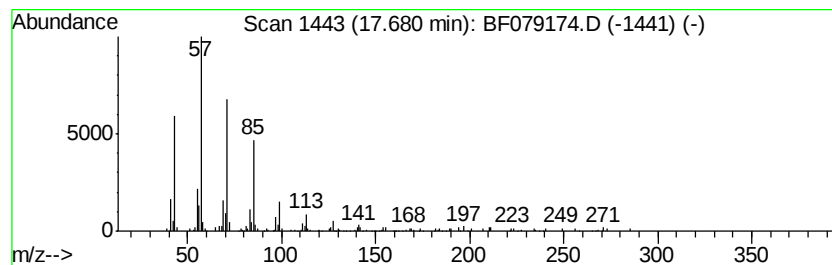
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Tetradecane, 2,6,10-trimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.68	6.22 ng	213260	Perylene-d12	17.95

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetradecane, 2,6,10-trimethyl-	240	C17H36	014905-56-7	87
2			Heptacosane	380	C27H56	000593-49-7	83
3			Nonadecane	268	C19H40	000629-92-5	83
4			Octadecane, 1-iodo-	380	C18H37I	000629-93-6	80
5			Tetracosane	338	C24H50	000646-31-1	80



Data Path : Z:\HPCHEM1\BNA F\DATA\BF051215\
Data File : BF079174.D
Acq On : 12 May 2015 23:06
Operator : TP/IZ
Sample : G2221-01
Misc :
ALS Vial : 19 Sample Multiplier: 1

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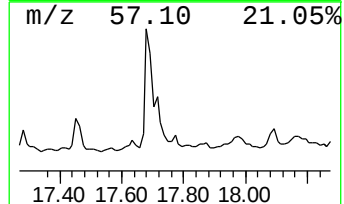
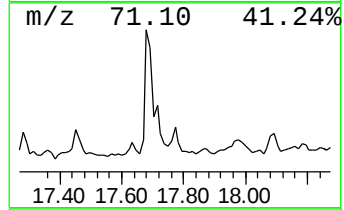
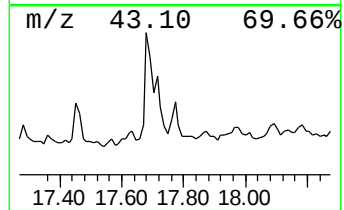
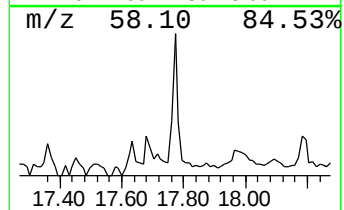
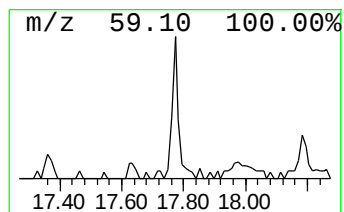
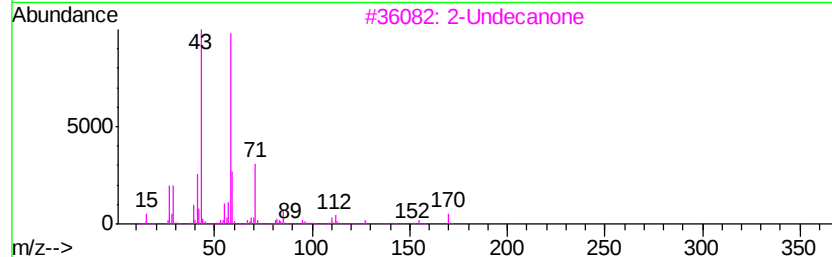
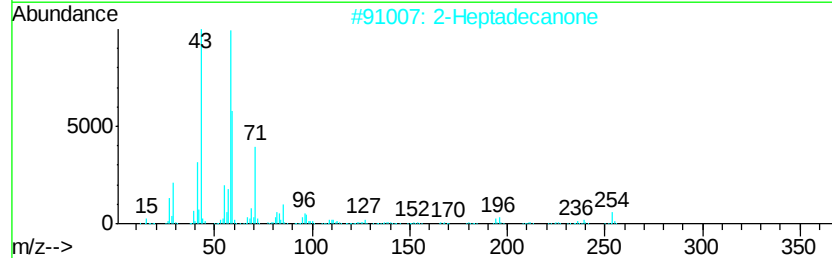
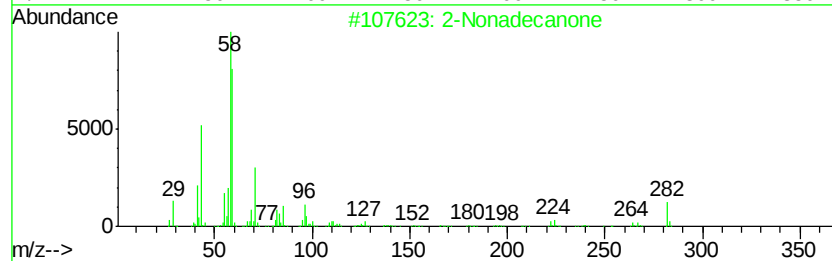
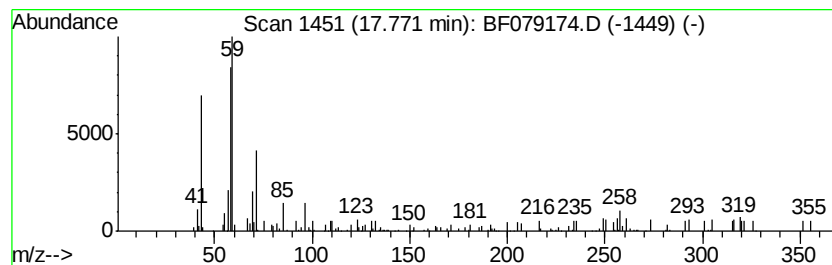
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 2-Nonadecanone Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.77	2.23 ng	76321	Perylene-d12	17.95

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Nonadecanone	282	C19H38O	000629-66-3	81
2			2-Heptadecanone	254	C17H34O	002922-51-2	52
3			2-Undecanone	170	C11H22O	000112-12-9	43
4			2-Pentadecanone, 6,10,14-trimethyl-	268	C18H36O	000502-69-2	37
5			2-Tetradecanone	212	C14H28O	002345-27-9	37



Data Path : Z:\HPCHEM1\BNA F\DATA\BF051215\
Data File : BF079174.D
Acq On : 12 May 2015 23:06
Operator : TP/IZ
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Misc :
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ClientSampled :
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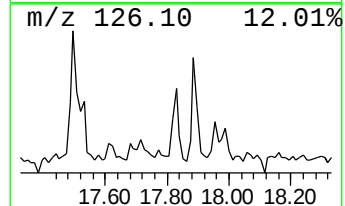
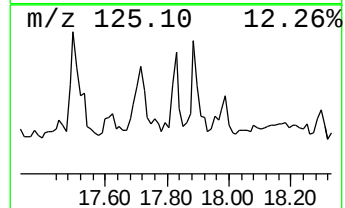
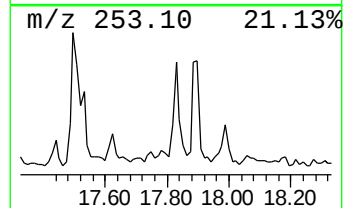
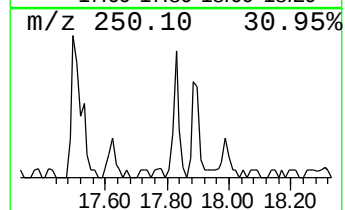
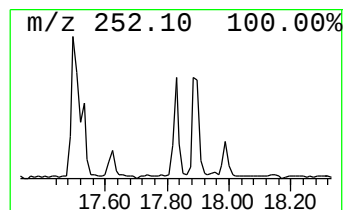
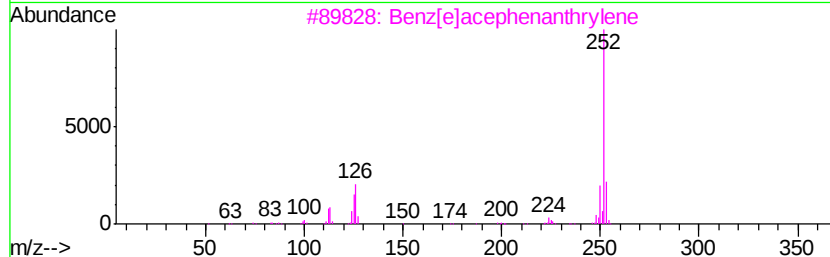
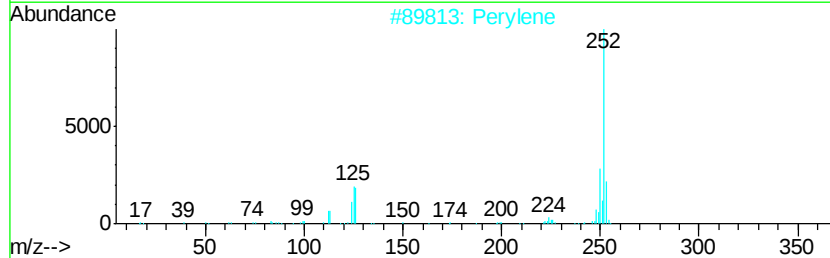
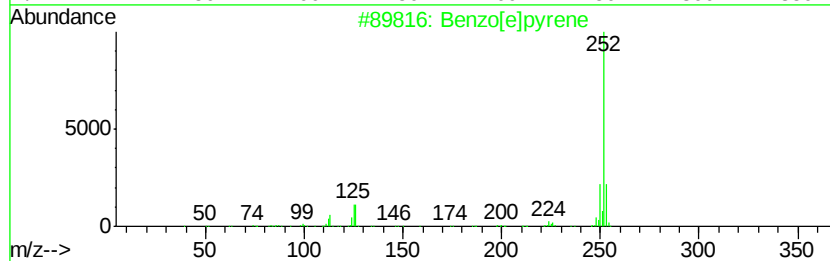
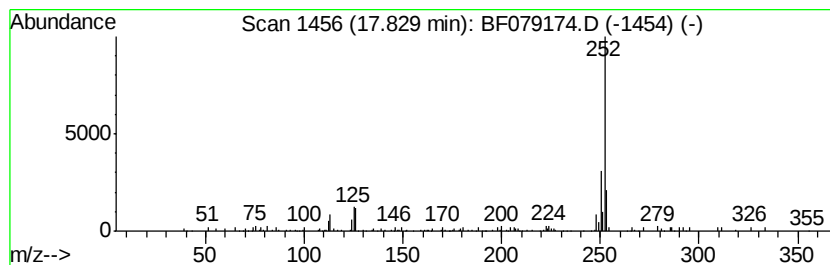
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Benzo[e]pyrene Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.83	2.14 ng	73315	Perylene-d12	17.95

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[e]pyrene	252	C20H12	000192-97-2	96
2			Perylene	252	C20H12	000198-55-0	93
3			Benzo[e]acephenanthrylene	252	C20H12	000205-99-2	91
4			Benzo[i]fluoranthene	252	C20H12	000205-82-3	90
5			Benzo[a]pyrene	252	C20H12	000050-32-8	90



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051215\
Data File : BF079174.D
Acq On : 12 May 2015 23:06
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Misc :
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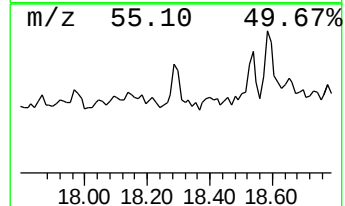
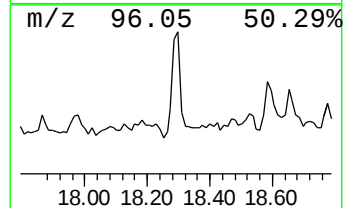
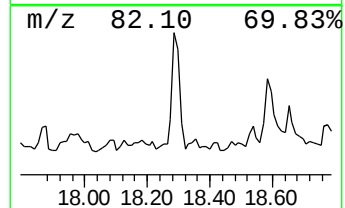
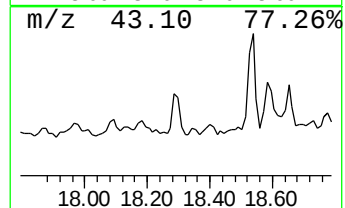
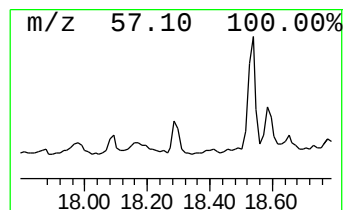
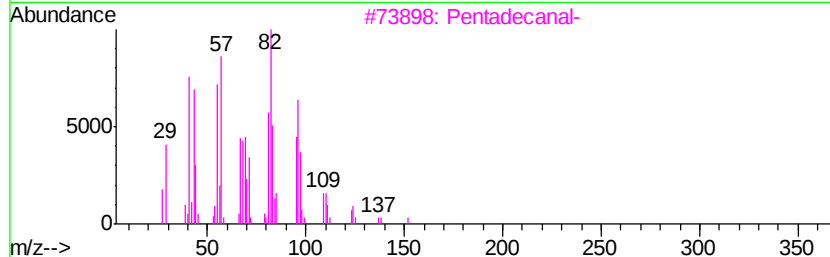
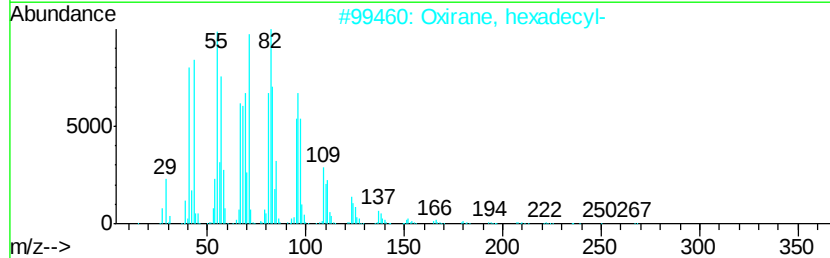
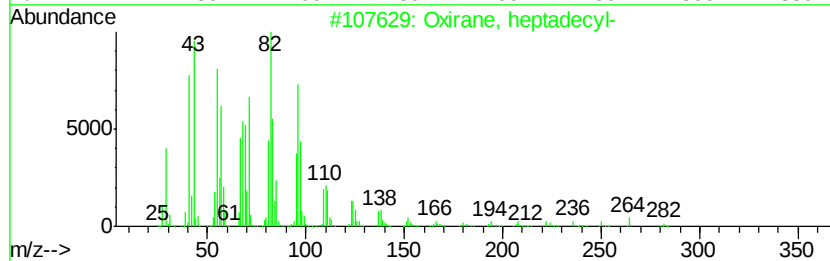
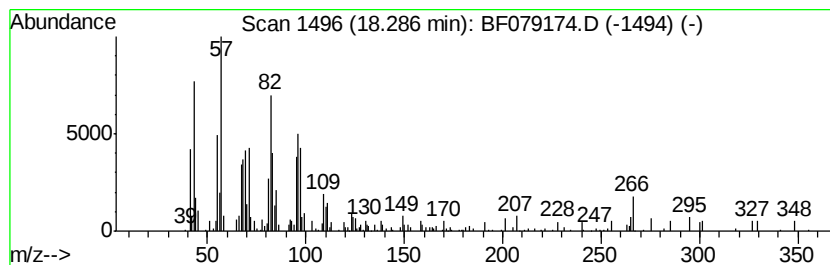
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Oxirane, heptadecyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.29	3.25 ng	111478	Perylene-d12	17.95

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Oxirane, heptadecyl-	282	C19H38O	067860-04-2	94
2		Oxirane, hexadecyl-	268	C18H36O	007390-81-0	90
3		Pentadecanal-	226	C15H30O	002765-11-9	64
4		Octadecanal	268	C18H36O	000638-66-4	58
5		Tetradecanal	212	C14H28O	000124-25-4	53



Data Path : Z:\HPCHEM1\BNA F\DATA\BF051215\
Data File : BF079174.D
Acq On : 12 May 2015 23:06
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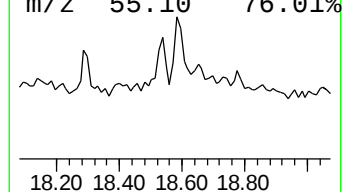
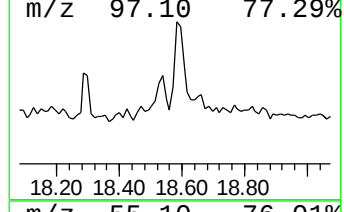
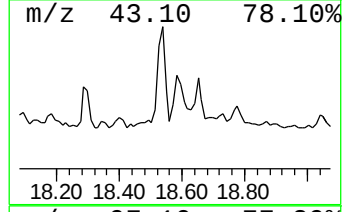
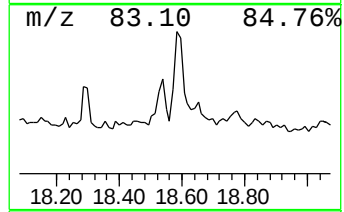
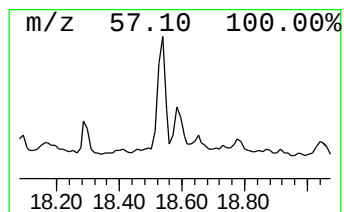
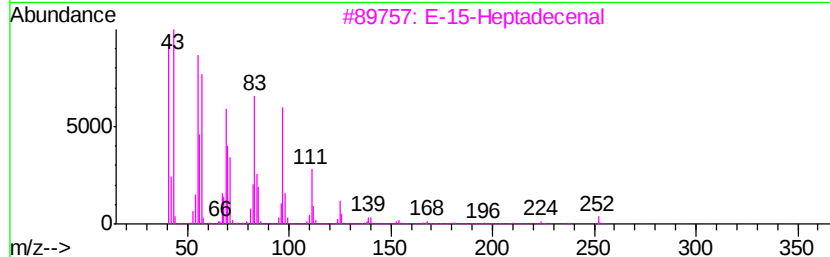
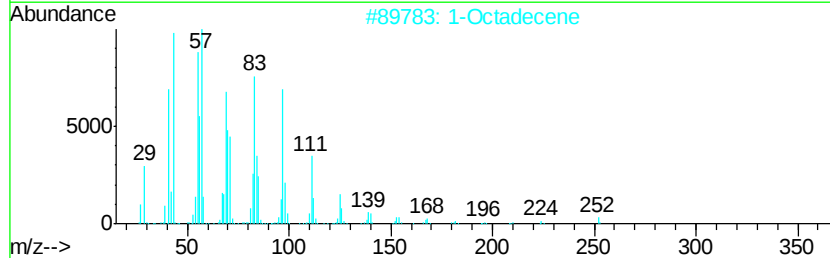
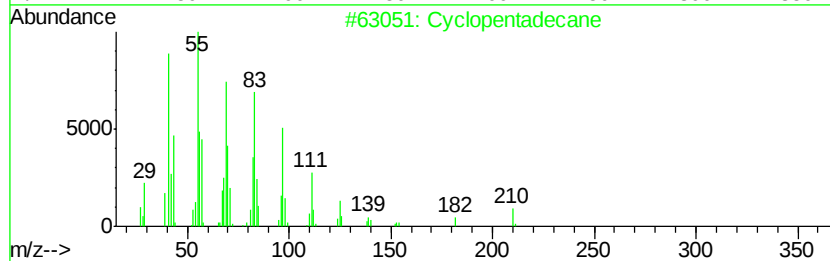
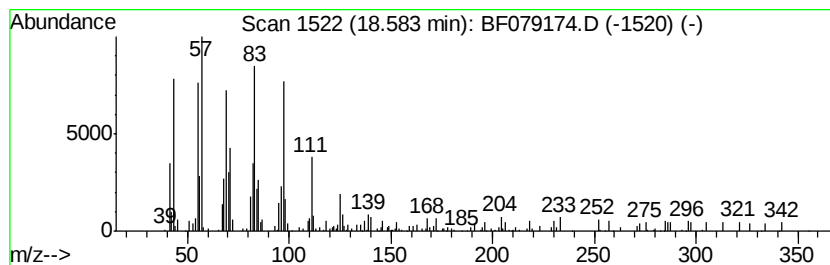
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Cyclopentadecane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.58	3.85 ng	132003	Perylene-d12	17.95

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentadecane	210	C15H30	000295-48-7	95
2		1-Octadecene	252	C18H36	000112-88-9	91
3		E-15-Heptadecenal	252	C17H32O	1000130-97-9	91
4		13-Tetradecen-1-ol acetate	254	C16H30O2	056221-91-1	91
5		1-Docosene	308	C22H44	001599-67-3	87



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051215\
Data File : BF079174.D
Acq On : 12 May 2015 23:06
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ClientSampleId :
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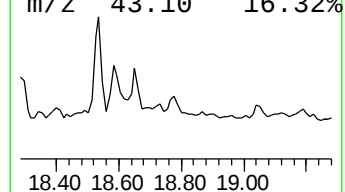
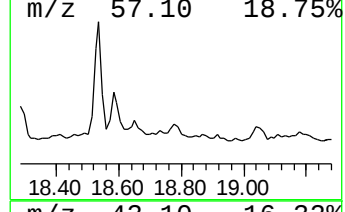
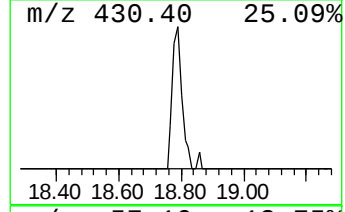
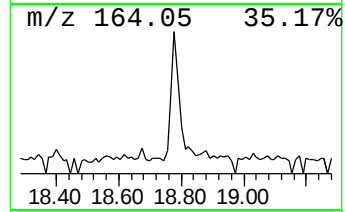
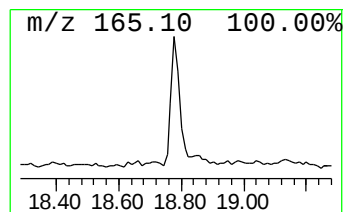
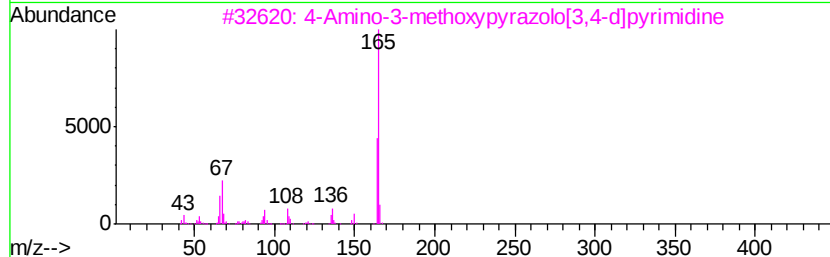
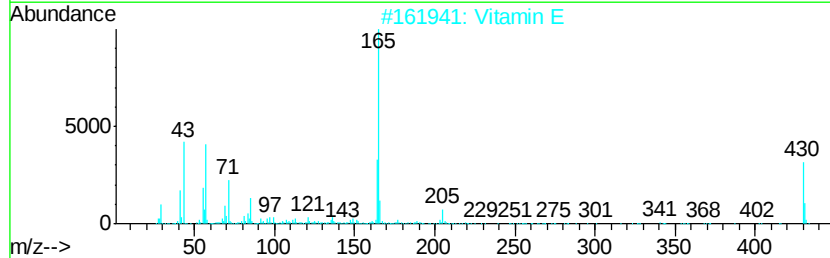
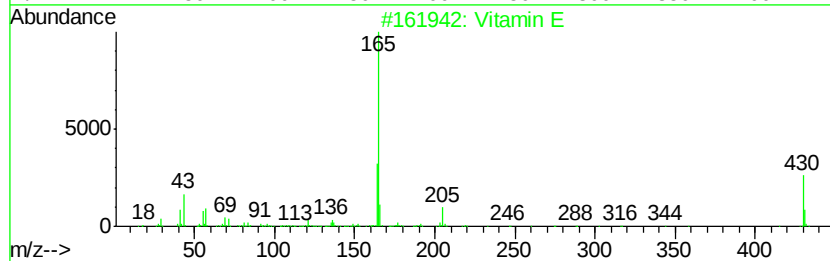
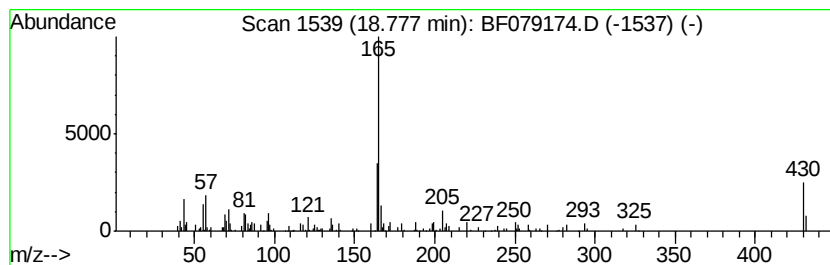
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 Vitamin E Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.78	2.46 ng	84359	Perylene-d12	17.95

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Vitamin E	430	C29H50O2	010191-41-0	94
2		Vitamin E	430	C29H50O2	000059-02-9	87
3		4-Amino-3-methoxypprazolo[3,4-d]...	165	C6H7N5O	111375-22-5	49
4		2,5-Pyridinedicarboxylic acid, d...	251	C13H17N04	000136-45-8	47
5		7-Methyl-8-oxo-1,2,3,4-tetrahydr...	165	C8H11N3O	026955-10-2	47



Data Path : Z:\HPCHEM1\BNA F\DATA\BF051215\
Data File : BF079174.D
Acq On : 12 May 2015 23:06
Operator : TP/IZ
Sample : G2221-01
Misc :
ALS Vial : 19 Sample Multiplier: 1

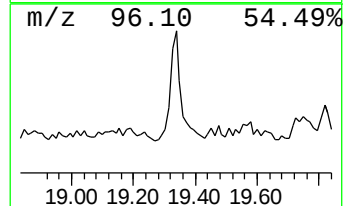
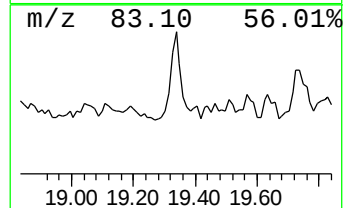
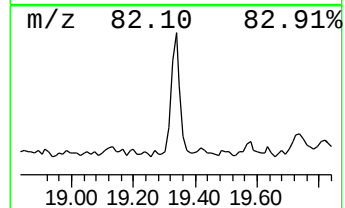
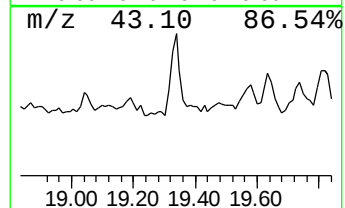
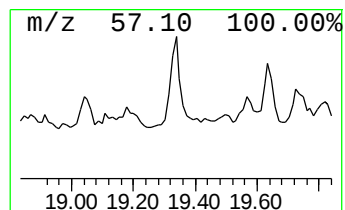
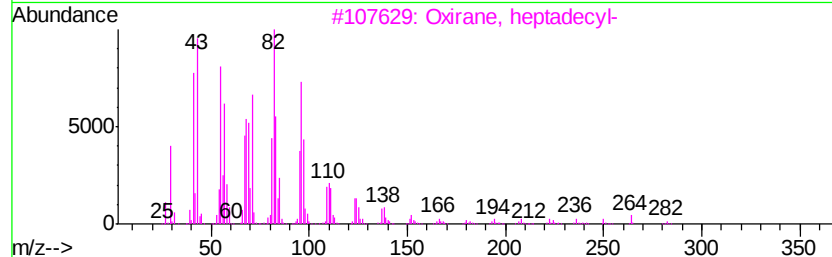
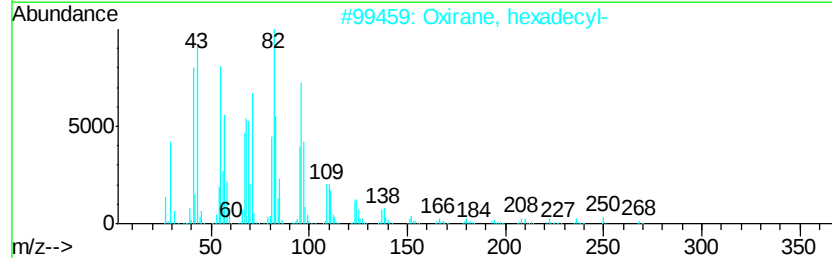
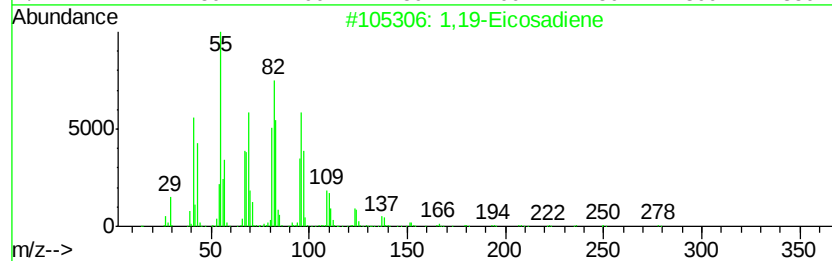
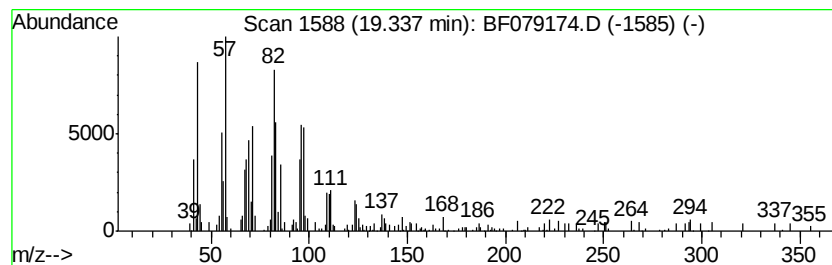
Instrument :
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ClientSampleId :
ALMASI-SOIL

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 18 1,19-Eicosadiene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.34	5.52 ng	189086	Perylene-d12	17.95	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,19-Eicosadiene	278	C20H38	014811-95-1	93
2	Oxirane, hexadecyl-	268	C18H36O	007390-81-0	87
3	Oxirane, heptadecyl-	282	C19H38O	067860-04-2	87
4	16-Heptadecenal	252	C17H32O	1000144-57-9	76
5	Octadecanal	268	C18H36O	000638-66-4	72



Data Path : Z:\HPCHEM1\BNA F\DATA\BF051215\
Data File : BF079174.D
Acq On : 12 May 2015 23:06
Operator : TP/IZ
Sample : G2221-01
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampled :
ALMASI-SOIL

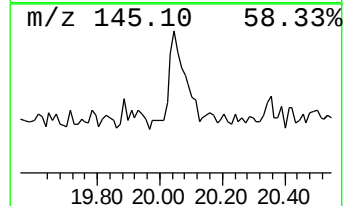
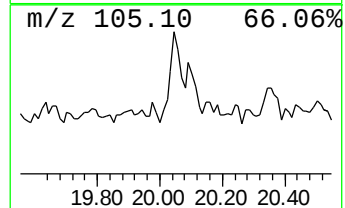
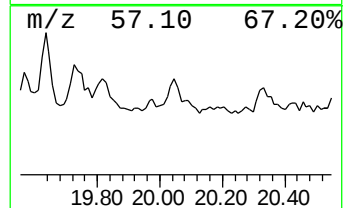
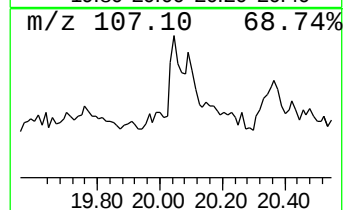
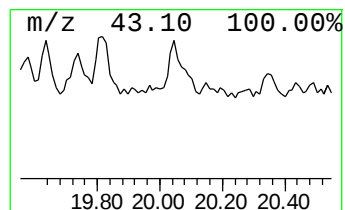
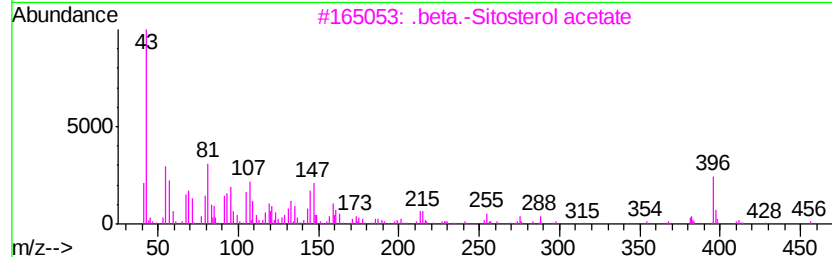
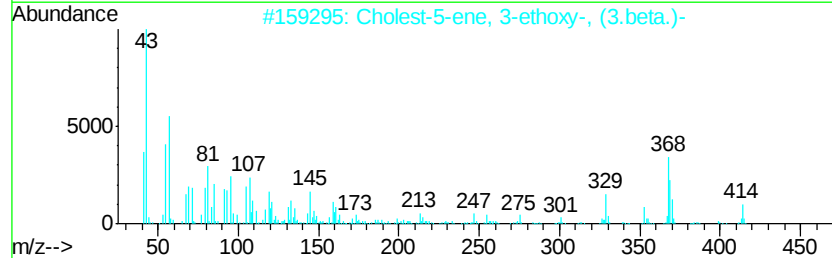
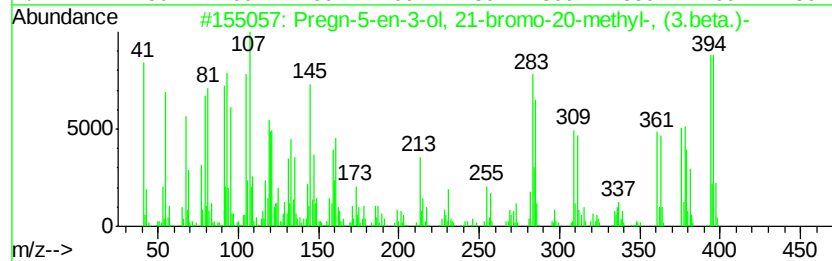
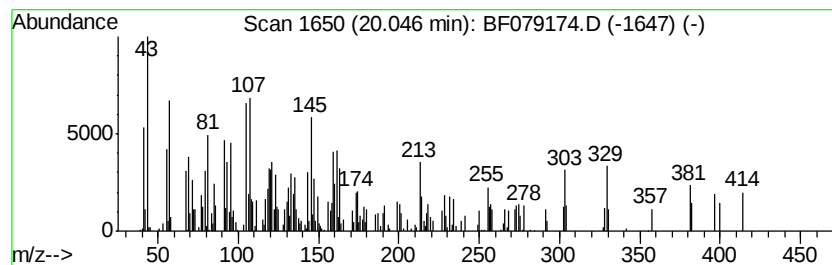
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF043015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 unknown20.05 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.05	4.55 ng	155984	Perylene-d12	17.95

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pregn-5-en-3-ol, 21-bromo-20-met...	394	C22H35BrO	055103-80-5	38
2		Cholest-5-ene, 3-ethoxy-, (3.beta...	414	C29H50O	000986-19-6	27
3		.beta.-Sitosterol acetate	456	C31H52O2	000915-05-9	25
4		Bisnor-7-desoxycholic acid	364	C22H36O4	1000252-00-7	25
5		Kauren-18-ol, acetate, (4.beta.)-	330	C22H34O2	072150-74-4	15



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051215\
 Data File : BF079174.D
 Acq On : 12 May 2015 23:06
 Operator : TP/IZ
 Sample : G2221-01
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 ALMASI-SOIL

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
unknown1.44	1.44	304.0	ng	4778730	1	7.28	314375	20.0
2-Pentanone, 4-hy...	5.07	9.8	ng	154124	1	7.28	314375	20.0
unknown6.99	6.99	87.2	ng	1370410	1	7.28	314375	20.0
unknown15.63	15.63	2.4	ng	82848	5	16.17	705429	20.0
Trichloroacetic a...	16.05	13.2	ng	467020	5	16.17	705429	20.0
1-Hexadecanesulfo...	16.88	10.7	ng	377707	5	16.17	705429	20.0
Tetradecane, 2,6,...	17.68	6.2	ng	213260	6	17.95	685416	20.0
2-Nonadecanone	17.77	2.2	ng	76321	6	17.95	685416	20.0
Benzo[e]pyrene	17.83	2.1	ng	73315	6	17.95	685416	20.0
Oxirane, heptadecyl-	18.29	3.3	ng	111478	6	17.95	685416	20.0
Cyclopentadecane	18.58	3.9	ng	132003	6	17.95	685416	20.0
Vitamin E	18.78	2.5	ng	84359	6	17.95	685416	20.0
1,19-Eicosadiene	19.34	5.5	ng	189086	6	17.95	685416	20.0
unknown20.05	20.05	4.5	ng	155984	6	17.95	685416	20.0