

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF071216\
 Data File : BF088940.D
 Acq On : 13 Jul 2016 1:24
 Operator : UM/SJ
 Sample : H3935-27
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 C3-01

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-BF063016.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	1.667	6	7	16	rVB	38373	86462	4.05%	0.498%
2	1.793	16	18	24	rVV	451986	594809	27.83%	3.429%
3	1.884	24	26	41	rVB	42217	92681	4.34%	0.534%
4	2.856	110	111	121	rBV2	29920	51129	2.39%	0.295%
5	4.856	283	286	291	rBV	161048	233569	10.93%	1.347%
6	5.302	322	325	328	rBV	1571976	1592492	74.51%	9.181%
7	6.353	414	417	420	rVB	1232734	1872295	87.60%	10.794%
8	6.479	425	428	430	rBV	1914446	2020717	94.54%	11.650%
9	6.707	446	448	449	rBV	367969	400693	18.75%	2.310%
10	6.856	459	461	464	rVB	1100381	1376060	64.38%	7.933%
11	7.268	494	497	500	rVB	884115	1176079	55.02%	6.781%
12	7.988	558	560	563	rVB	630267	532877	24.93%	3.072%
13	9.073	652	655	657	rBV	2044615	2137377	100.00%	12.323%
14	9.462	687	689	691	rBV	336369	275024	12.87%	1.586%
15	9.553	691	697	700	rVB2	15028	33314	1.56%	0.192%
16	9.691	707	709	711	rVB	22106	21848	1.02%	0.126%
17	9.736	711	713	716	rVB	582468	531963	24.89%	3.067%
18	10.182	751	752	754	rVB	36172	28860	1.35%	0.166%
19	10.525	780	782	784	rBV	1093971	992396	46.43%	5.722%
20	10.651	792	793	797	rBV2	35533	64011	2.99%	0.369%
21	10.971	817	821	824	rBV2	70958	113235	5.30%	0.653%
22	11.096	831	832	834	rBV	52960	63004	2.95%	0.363%
23	11.211	840	842	845	rVB	440270	404192	18.91%	2.330%
24	11.279	846	848	853	rVB2	13794	29041	1.36%	0.167%
25	11.451	861	863	868	rVV3	29919	42927	2.01%	0.247%
26	11.531	868	870	873	rVB	24953	31659	1.48%	0.183%
27	11.759	888	890	892	rVV	55214	41334	1.93%	0.238%
28	11.817	892	895	901	rVB	36861	59855	2.80%	0.345%
29	12.800	978	981	983	rBV	1365117	1212255	56.72%	6.989%
30	13.154	1010	1012	1015	rBV3	12711	22294	1.04%	0.129%
31	13.382	1026	1032	1035	rBV6	13275	33398	1.56%	0.193%
32	13.702	1057	1060	1064	rBV	133227	148904	6.97%	0.858%
33	13.840	1069	1072	1074	rBV	428342	413115	19.33%	2.382%
34	14.331	1113	1115	1121	rBV	77587	110892	5.19%	0.639%

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35	14.685	1144	1146	1148	rVB	21744	22577	1.06%	0.130%
36	14.834	1157	1159	1162	rBV3	14455	24779	1.16%	0.143%
37	15.211	1189	1192	1195	rBV	294319	336438	15.74%	1.940%
38	15.691	1232	1234	1238	rVB4	17204	31884	1.49%	0.184%
39	16.651	1316	1318	1322	rVB4	18398	40633	1.90%	0.234%
40	17.451	1385	1388	1391	rVB5	21328	47941	2.24%	0.276%

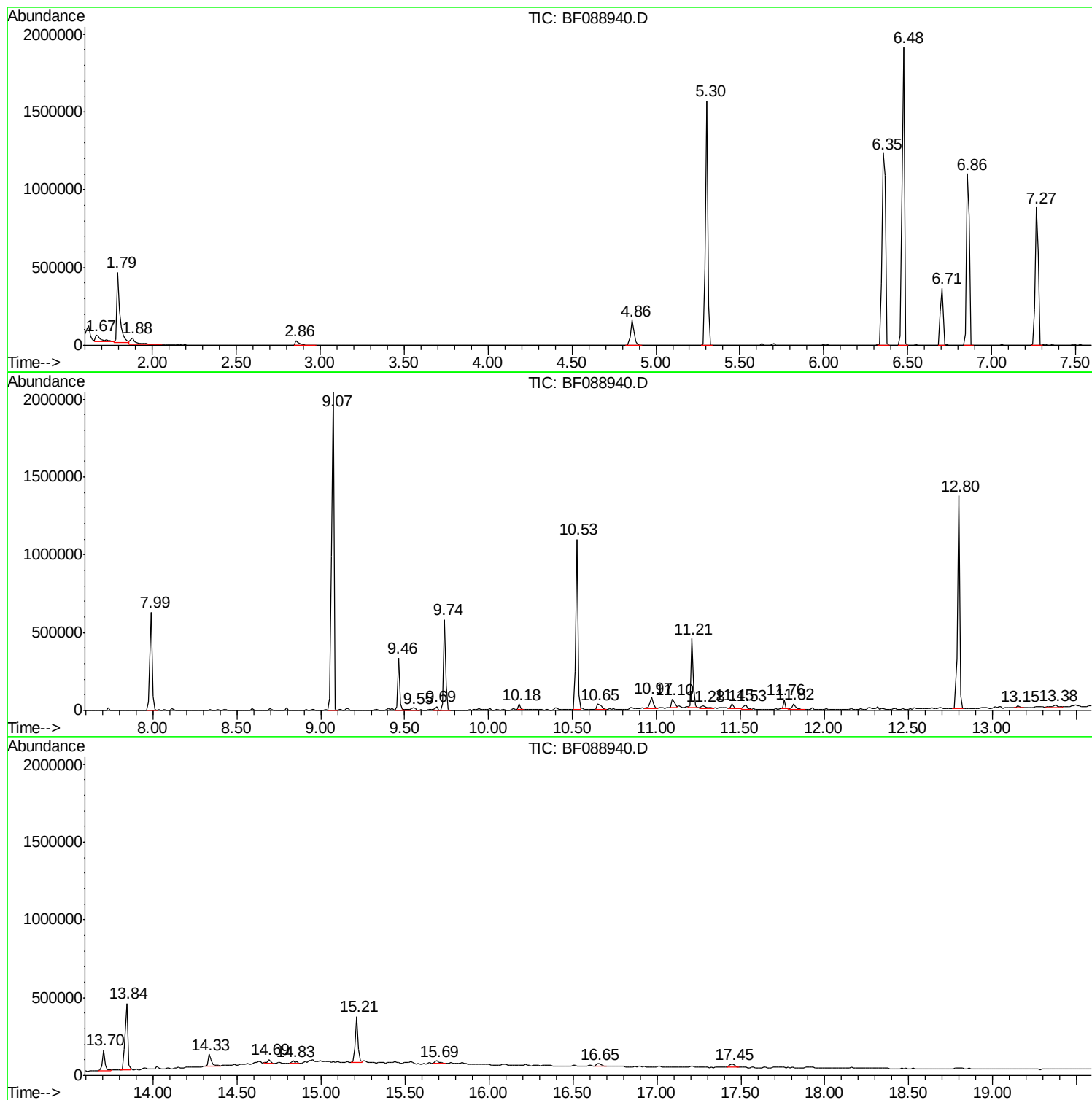
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF063016.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P



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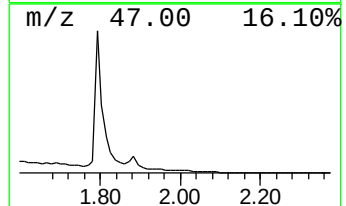
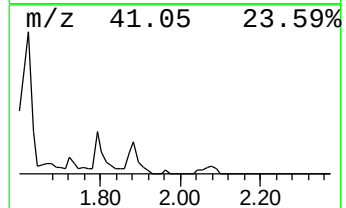
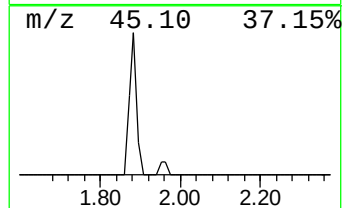
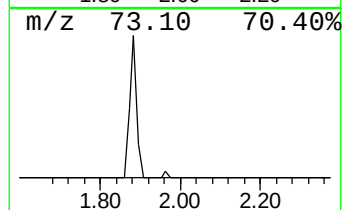
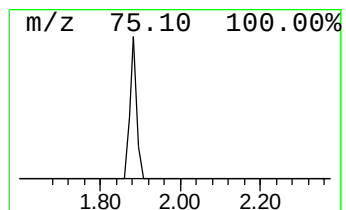
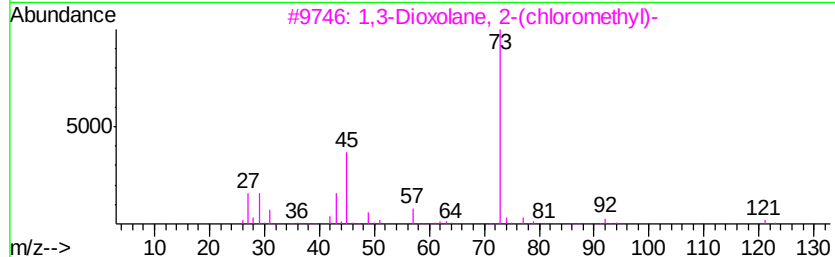
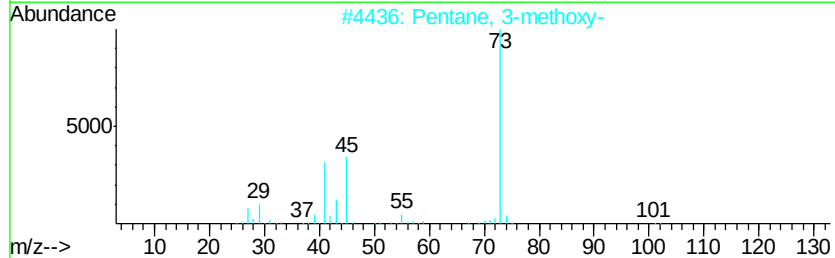
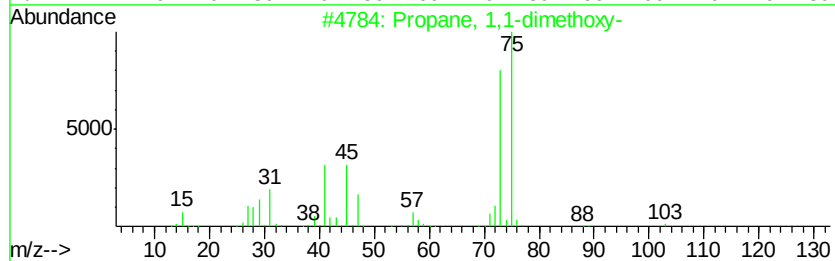
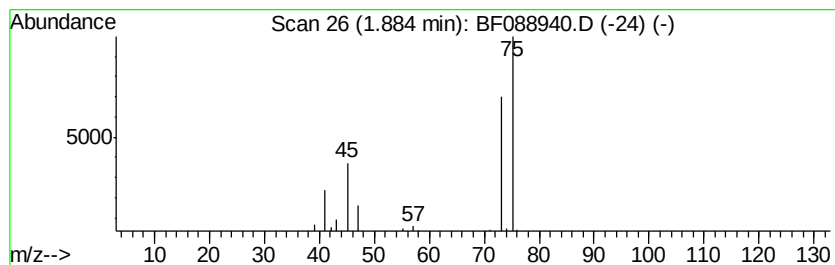
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TIC Library : C:\Database\NIST11.L
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Peak Number 3 Propane, 1,1-dimethoxy- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.88	4.63 ng	92681	1,4-Dichlorobenzene-d4	6.71

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Propane, 1,1-dimethoxy-	104	C5H12O2	004744-10-9	78
2		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	9
3		1,3-Dioxolane, 2-(chloromethyl)-	122	C4H7ClO2	002568-30-1	9
4		Glycine	75	C2H5NO2	000056-40-6	7
5		1-Propene-1-thiol	74	C3H6S	000925-89-3	7



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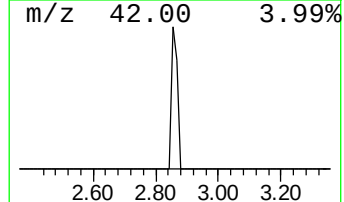
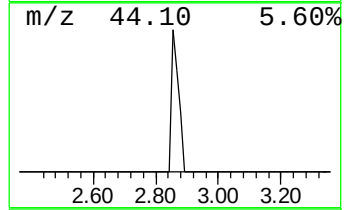
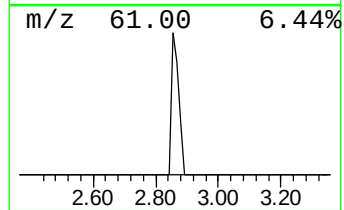
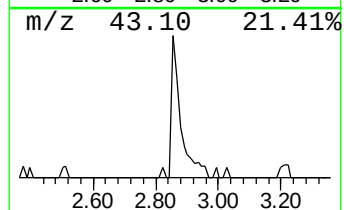
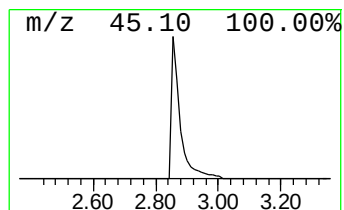
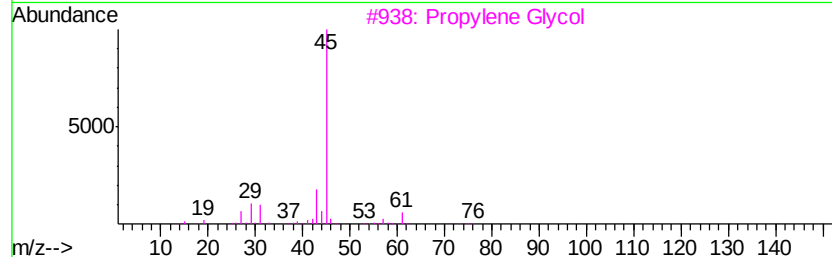
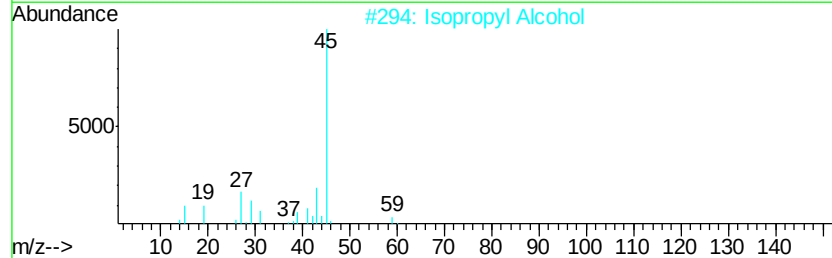
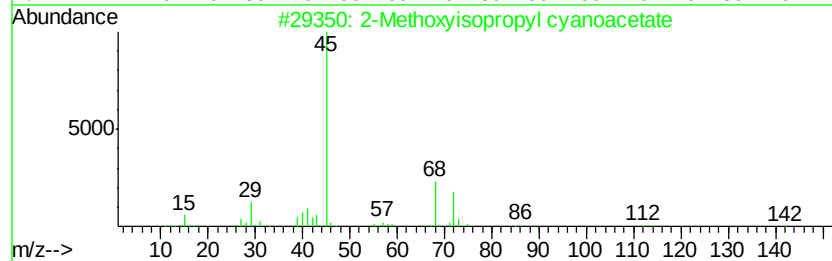
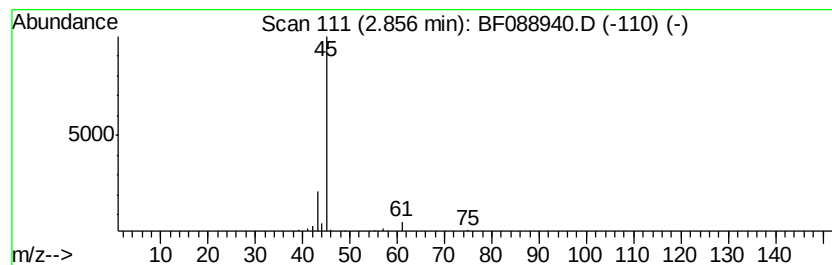
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TIC Library : C:\Database\NIST11.L
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Peak Number 4 unknown2.86 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.86	2.55 ng	51129	1,4-Dichlorobenzene-d4	6.71

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Methoxyisopropyl cyanoacetate	157	C7H11NO3	032804-79-8	9
2		Isopropyl Alcohol	60	C3H8O	000067-63-0	9
3		Propylene Glycol	76	C3H8O2	000057-55-6	9
4		Methoxyacetic acid, pentyl ester	160	C8H16O3	168920-35-2	9
5		Propane, 2-ethoxy-	88	C5H12O	000625-54-7	9



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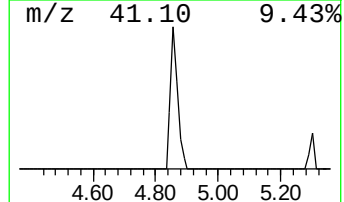
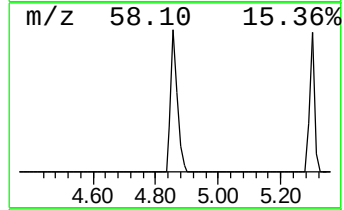
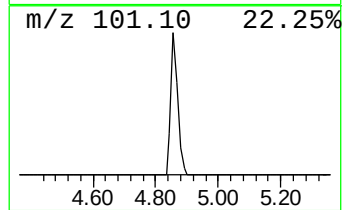
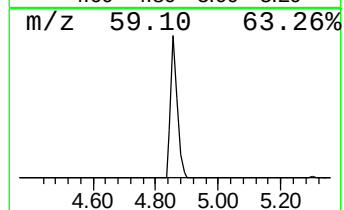
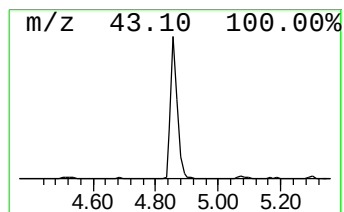
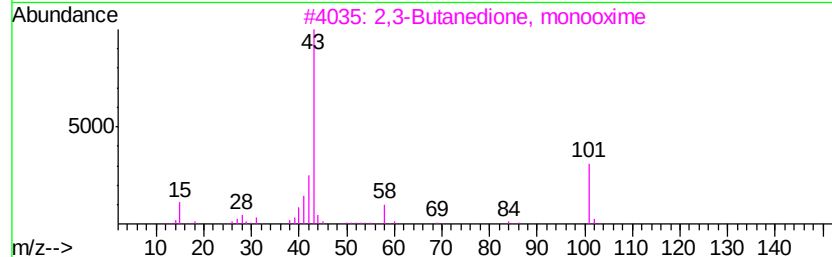
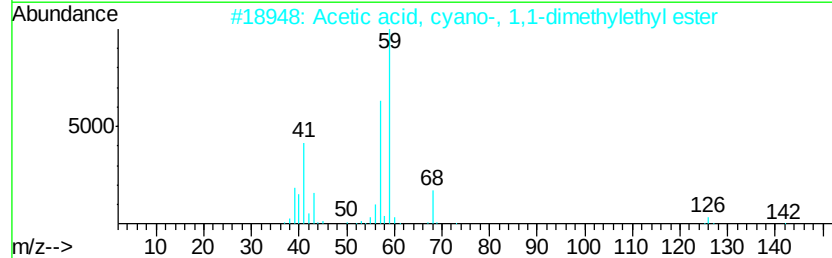
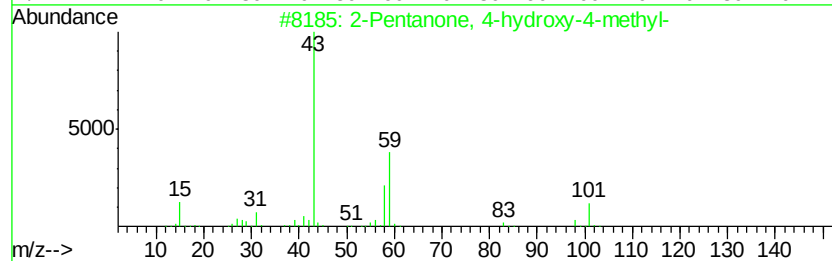
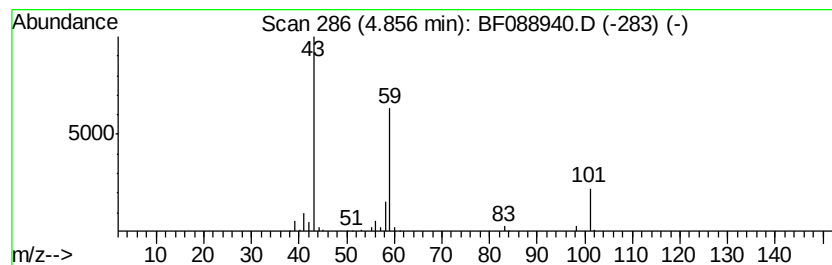
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TIC Library : C:\Database\NIST11.L
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Peak Number 5 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.86	11.66 ng	233569	1,4-Dichlorobenzene-d4	6.71

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56	
2	Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	17	
3	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	17	
4	Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9	
5	Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9	



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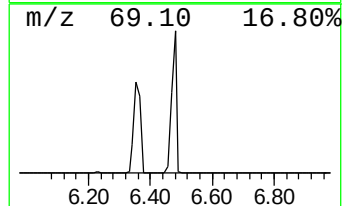
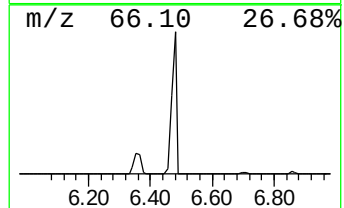
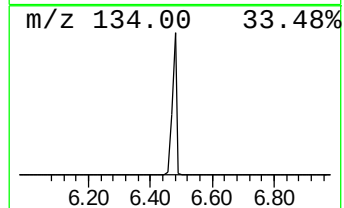
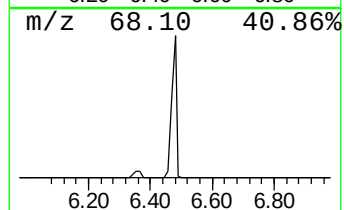
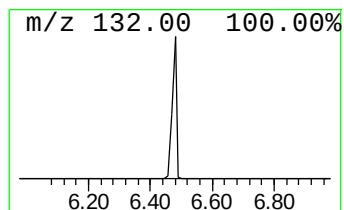
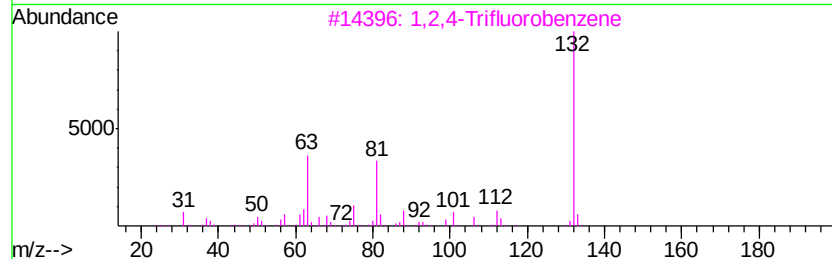
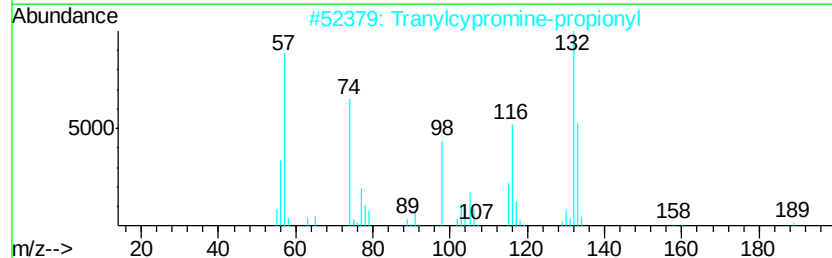
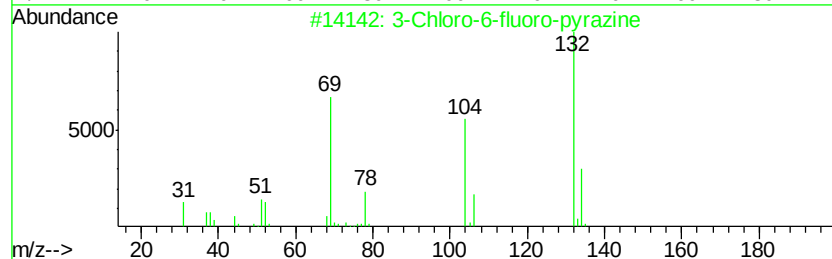
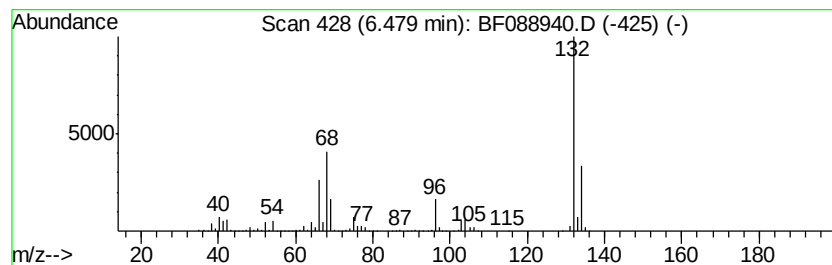
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TIC Library : C:\Database\NIST11.L
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Peak Number 6 unknown6.48 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.48	100.86 ng	2020720	1,4-Dichlorobenzene-d4	6.71

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3		1,2,4-Trifluorobenzene	132	C6H3F3	000367-23-7	9
4		1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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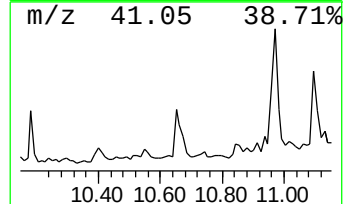
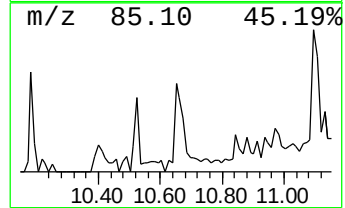
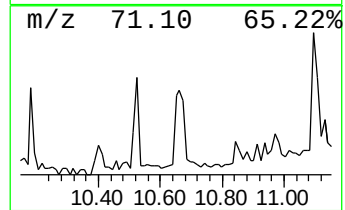
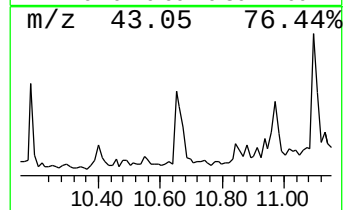
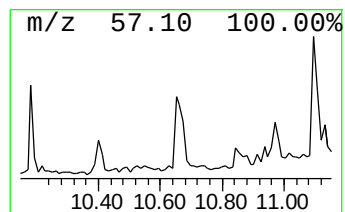
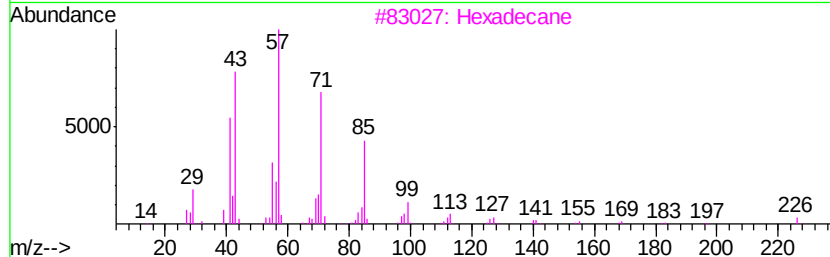
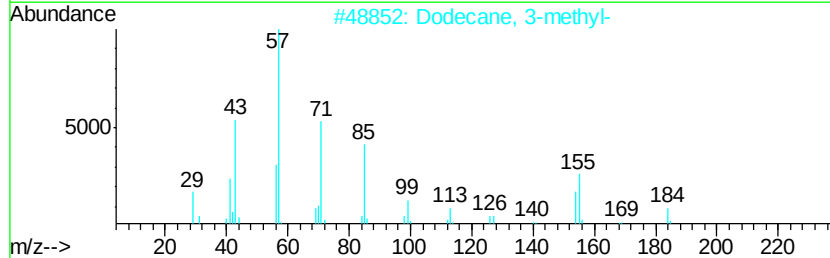
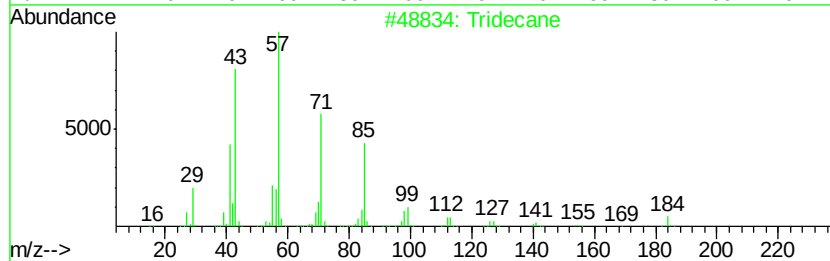
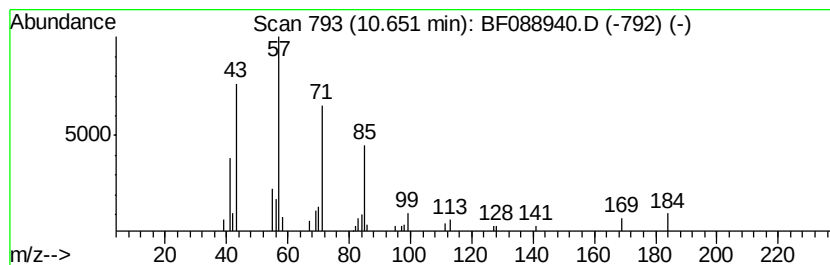
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Peak Number 8 Tridecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.65	3.17 ng	64011	Phenanthrene-d10	11.21

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tridecane	184	C13H28	000629-50-5	93
2		Dodecane, 3-methyl-	184	C13H28	017312-57-1	83
3		Hexadecane	226	C16H34	000544-76-3	83
4		Heneicosane	296	C21H44	000629-94-7	83
5		Octacosane	394	C28H58	000630-02-4	80



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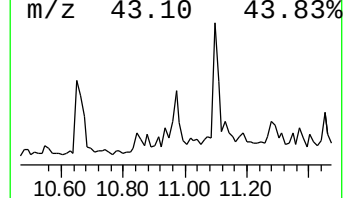
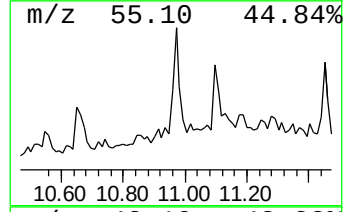
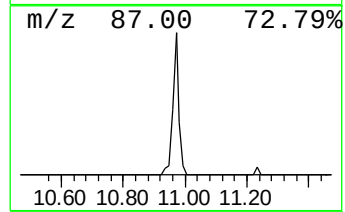
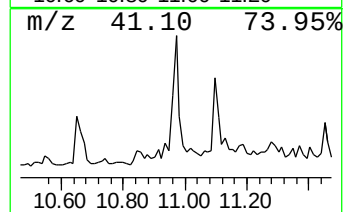
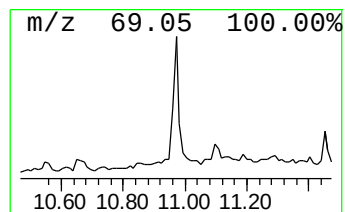
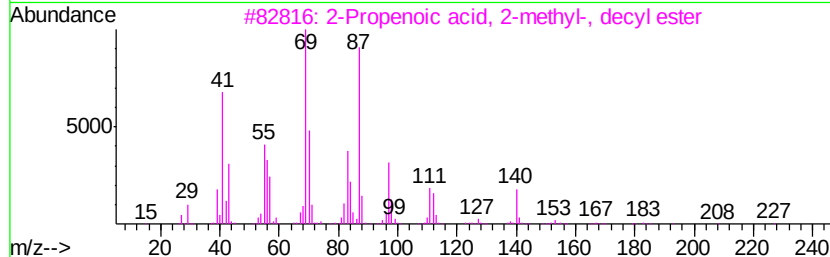
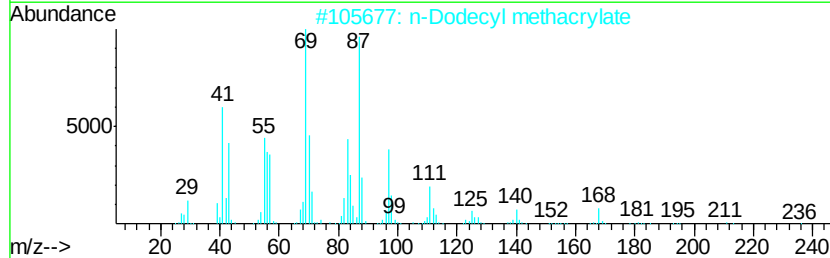
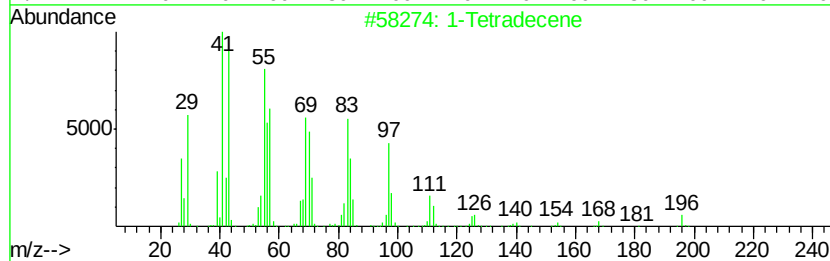
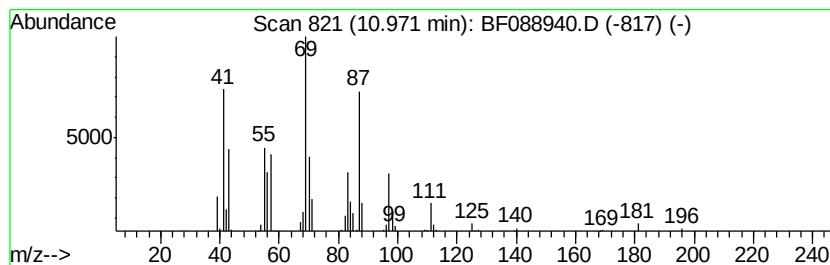
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ClientSampled :
C3-01

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF063016.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 1-Tetradecene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.97	5.60 ng	113235	Phenanthrene-d10	11.21	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Tetradecene	196	C14H28	001120-36-1	93
2	n-Dodecyl methacrylate	254	C16H30O2	000142-90-5	91
3	2-Propenoic acid, 2-methyl-, dec...	226	C14H26O2	003179-47-3	91
4	Methacrylic acid, tetradecyl ester	282	C18H34O2	1000340-29-0	60
5	Methacrylic acid, tridecyl ester	268	C17H32O2	1000340-28-9	53



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF071216\
Data File : BF088940.D
Acq On : 13 Jul 2016 1:24
Operator : UM/SJ
Sample : H3935-27
Misc :
ALS Vial : 19 Sample Multiplier: 1

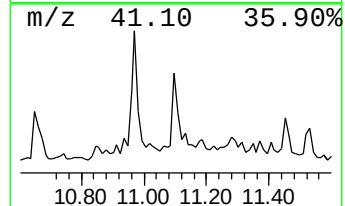
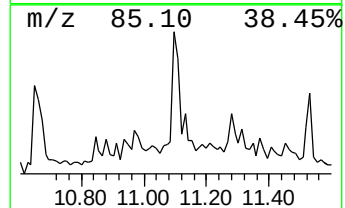
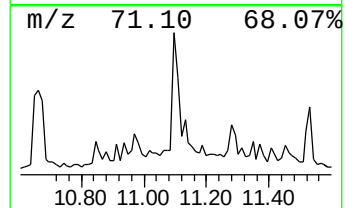
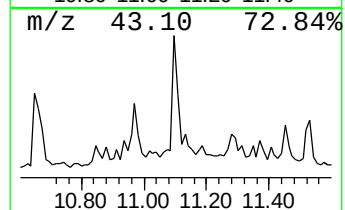
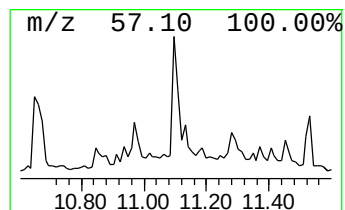
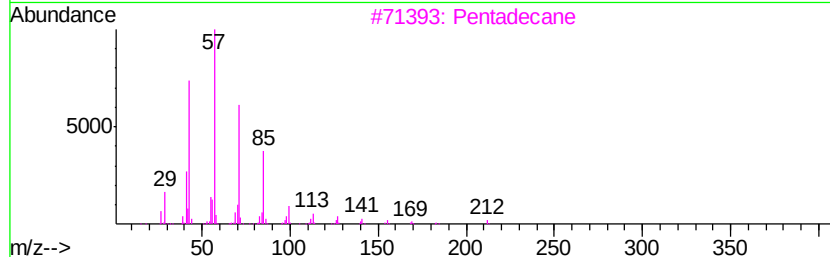
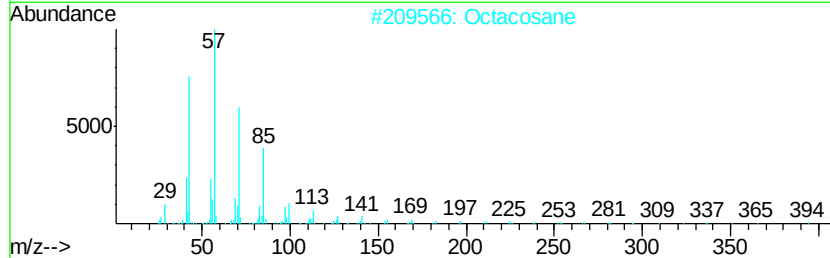
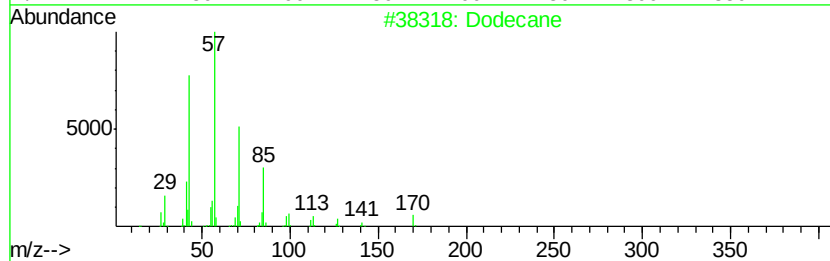
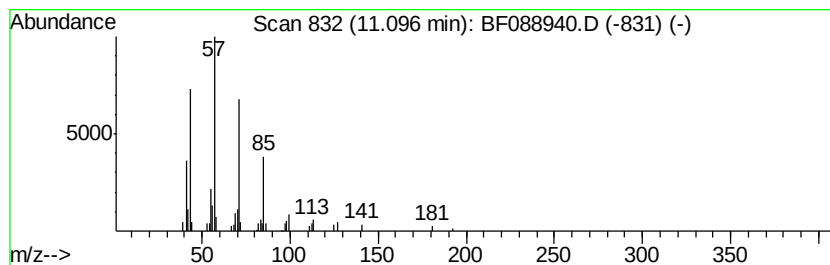
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Dodecane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD		R.T.
11.10	3.12 ng	63004	Phenanthrene-d10		11.21
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dodecane	170	C12H26	000112-40-3	90
2	Octacosane	394	C28H58	000630-02-4	90
3	Pentadecane	212	C15H32	000629-62-9	90
4	Eicosane	282	C20H42	000112-95-8	86
5	Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	86



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF071216\
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Acq On : 13 Jul 2016 1:24
Operator : UM/SJ
Sample : H3935-27
Misc :
ALS Vial : 19 Sample Multiplier: 1

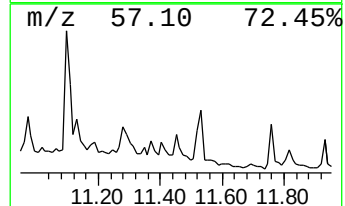
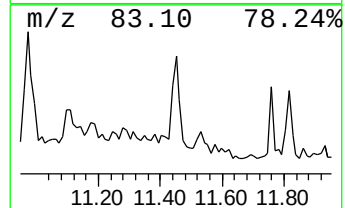
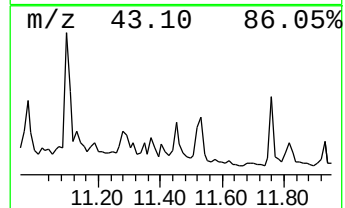
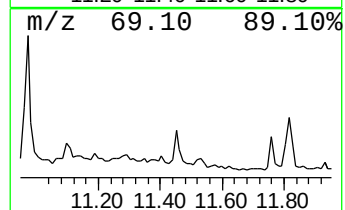
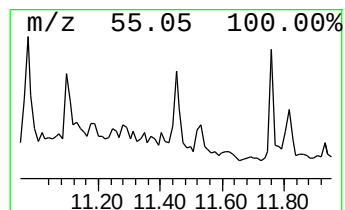
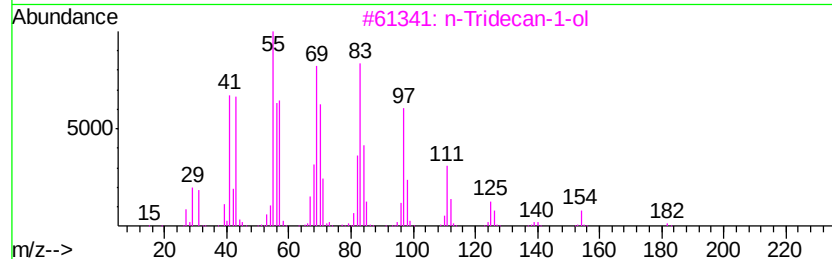
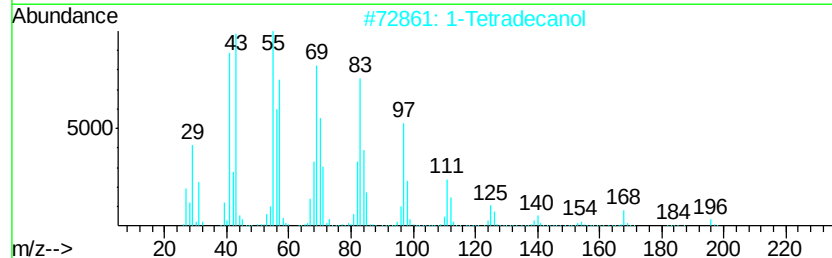
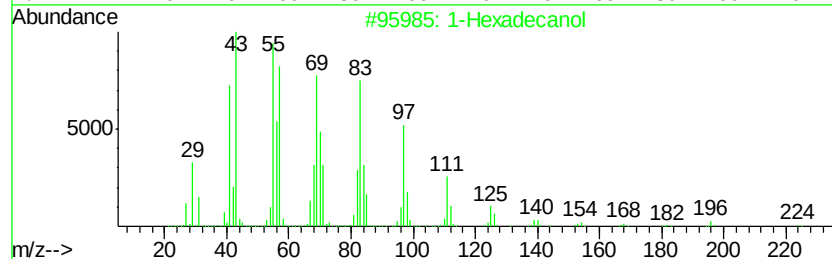
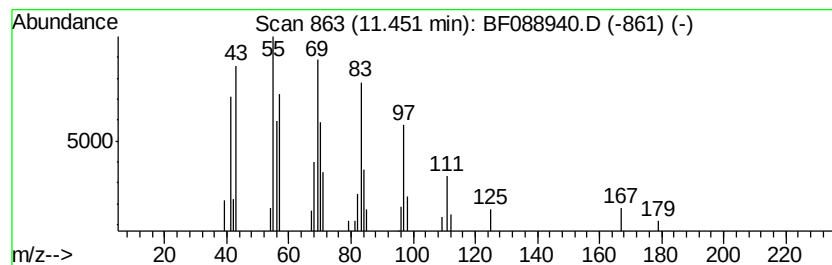
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ClientSampled :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF063016.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 1-Hexadecanol Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD		R.T.
11.45	2.12 ng	42927	Phenanthrene-d10		11.21
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexadecanol	242	C16H34O	036653-82-4	91
2	1-Tetradecanol	214	C14H30O	000112-72-1	91
3	n-Tridecan-1-ol	200	C13H28O	000112-70-9	91
4	9-Octadecene, (E)-	252	C18H36	007206-25-9	83
5	1-Dodecanol	186	C12H26O	000112-53-8	83



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF071216\
Data File : BF088940.D
Acq On : 13 Jul 2016 1:24
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Sample : H3935-27
Misc :
ALS Vial : 19 Sample Multiplier: 1

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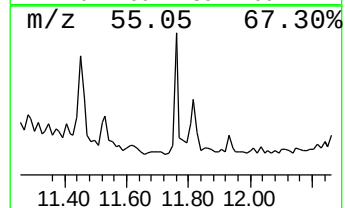
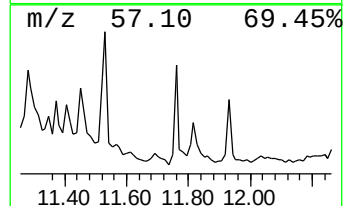
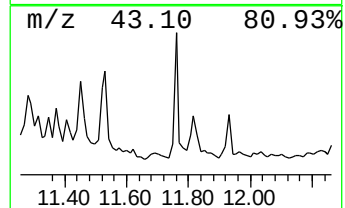
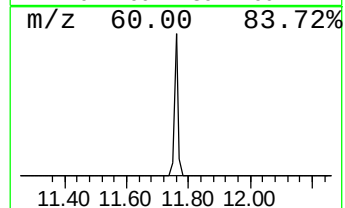
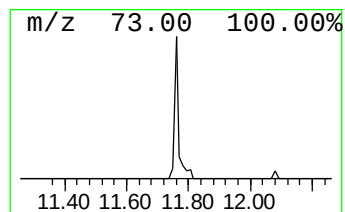
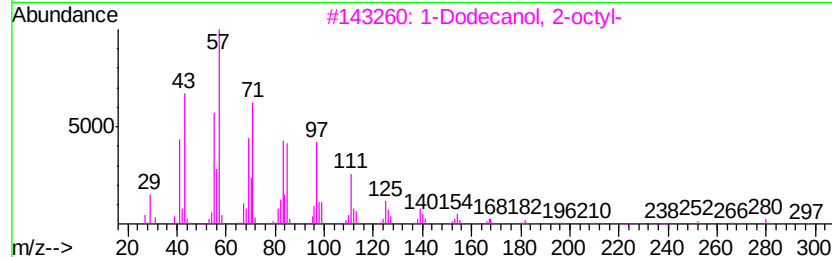
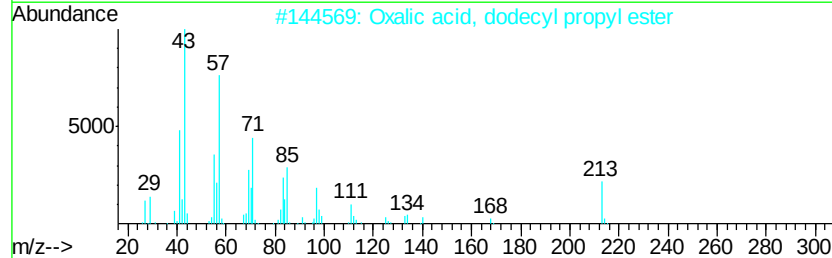
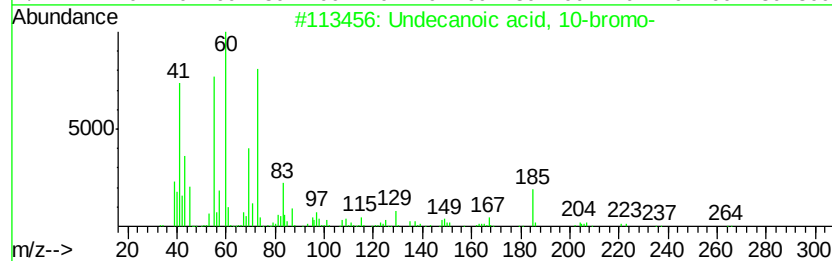
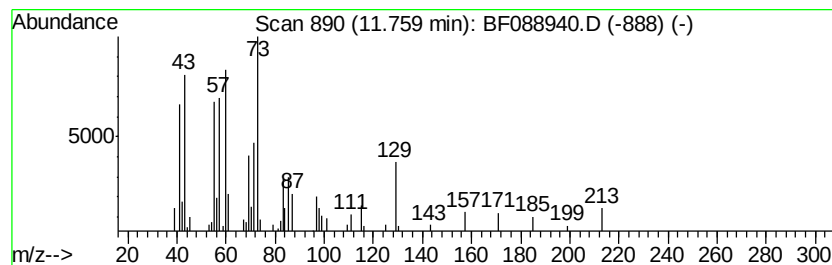
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 unknown11.76 Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.76	2.05 ng	41334	Phenanthrene-d10	11.21

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Undecanoic acid, 10-bromo-	264	C11H21BrO2	018294-93-4	38
2			Oxalic acid, dodecyl propyl ester	300	C17H32O4	1000309-26-5	22
3			1-Dodecanol, 2-octyl-	298	C20H42O	005333-42-6	18
4			Oxalic acid, allyl pentadecyl ester	340	C20H36O4	1000309-24-3	18
5			Oxalic acid, allyl hexadecyl ester	354	C21H38O4	1000309-24-4	15



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF071216\
Data File : BF088940.D
Acq On : 13 Jul 2016 1:24
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Sample : H3935-27
Misc :
ALS Vial : 19 Sample Multiplier: 1

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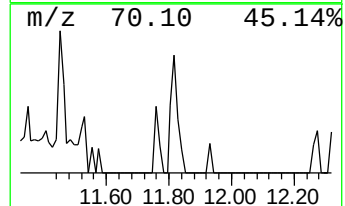
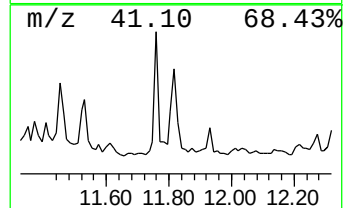
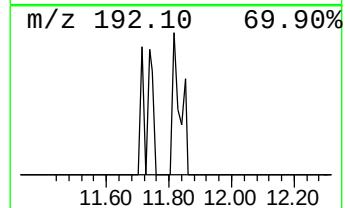
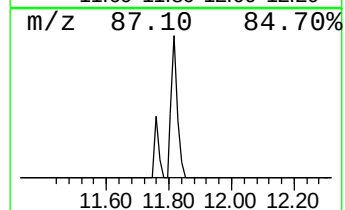
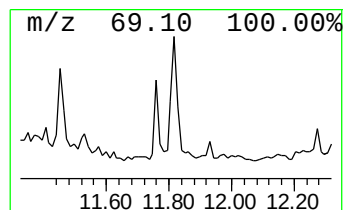
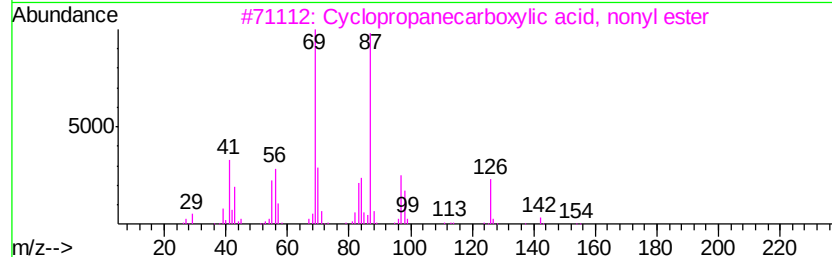
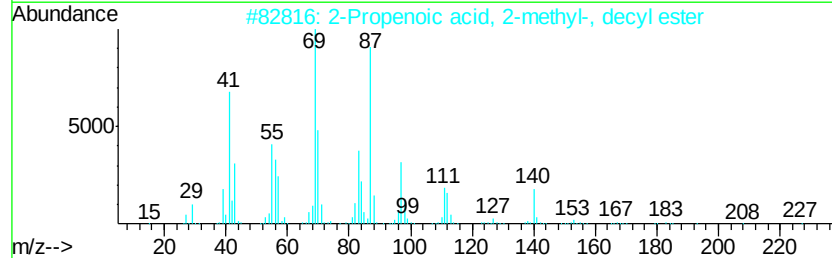
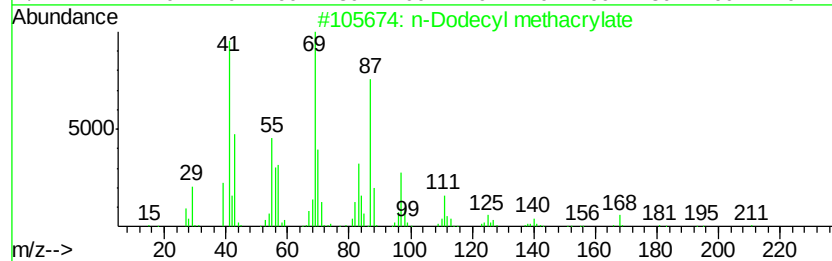
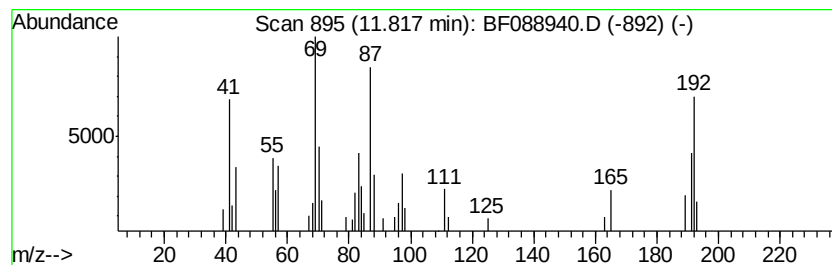
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 unknown11.82 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.82	2.96 ng	59855	Phenanthrene-d10	11.21

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Dodecyl methacrylate	254	C16H30O2	000142-90-5	49
2			2-Propenoic acid, 2-methyl-, dec...	226	C14H26O2	003179-47-3	47
3			Cyclopropanecarboxylic acid, non...	212	C13H24O2	060128-06-5	43
4			Methacrylic acid, tetradecyl ester	282	C18H34O2	1000340-29-0	41
5			Methacrylic acid, tridecyl ester	268	C17H32O2	1000340-28-9	38



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF071216\
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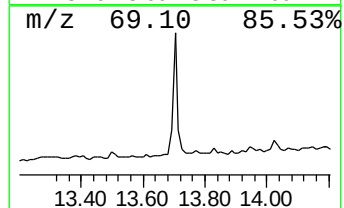
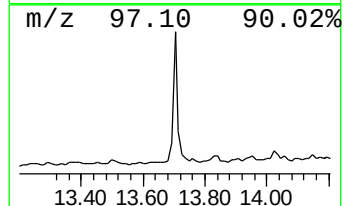
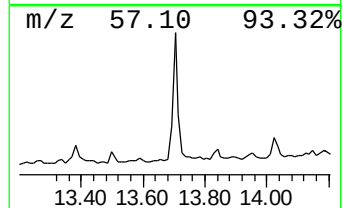
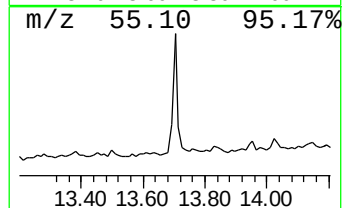
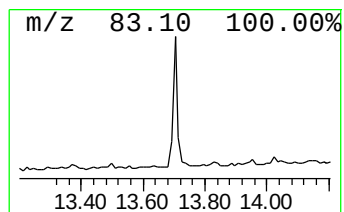
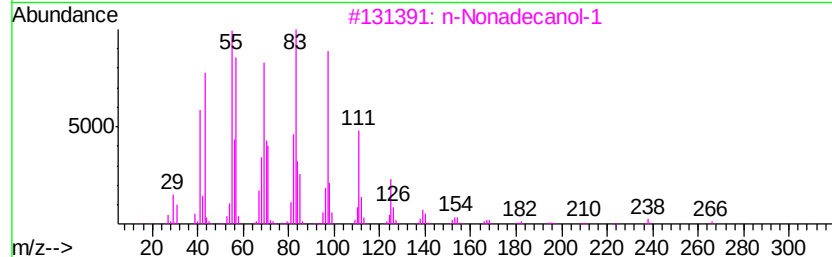
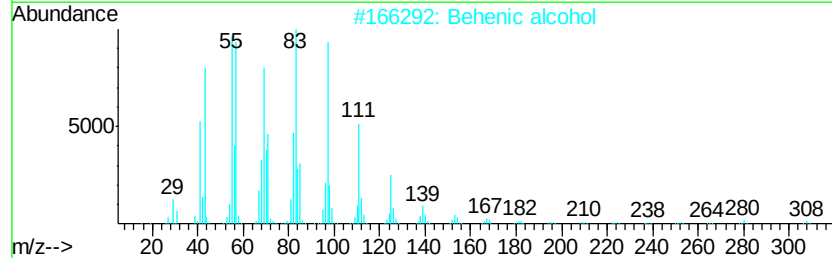
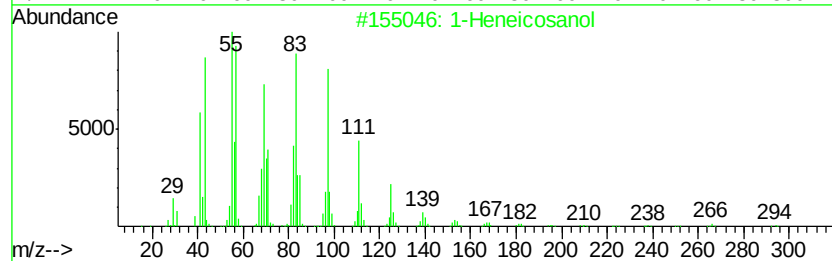
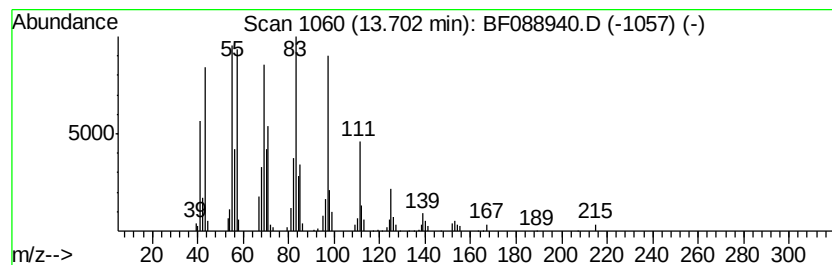
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 1-Heneicosanol Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.70	7.21 ng	148904	Chrysene-d12	13.84	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heneicosanol	312	C21H44O	015594-90-8	95
2	Behenic alcohol	326	C22H46O	000661-19-8	95
3	n-Nonadecanol-1	284	C19H40O	001454-84-8	95
4	Trifluoroacetoxv hexadecane	338	C18H33F3O2	006222-03-3	93
5	Pentafluoropropionic acid, penta...	374	C18H31F5O2	959092-08-1	93



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF071216\
Data File : BF088940.D
Acq On : 13 Jul 2016 1:24
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Misc :
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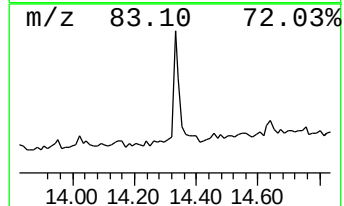
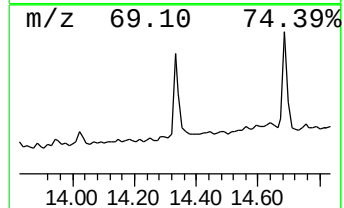
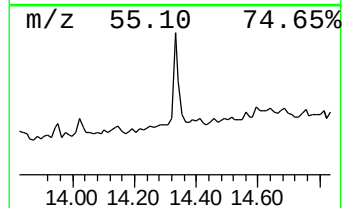
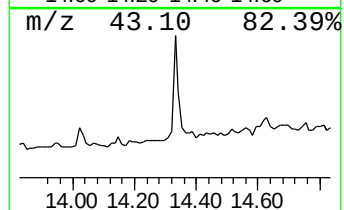
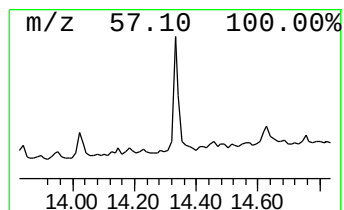
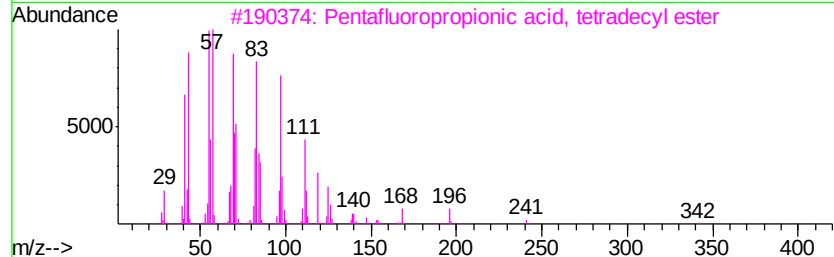
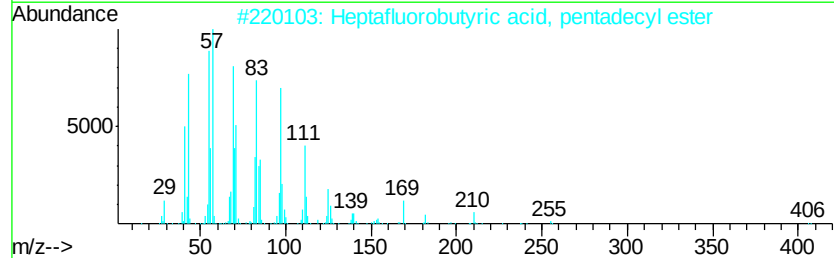
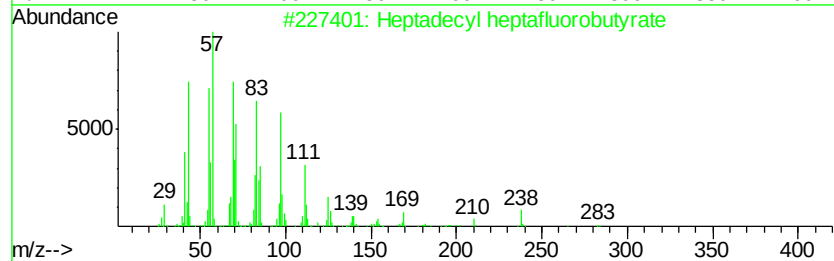
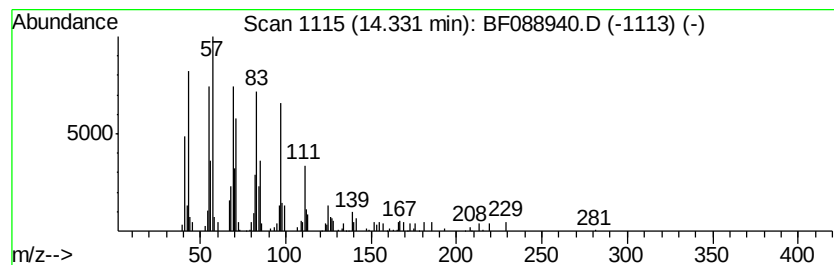
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 15 Heptadecyl heptafluorobutyrate Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.33	5.37 ng	110892	Chrysene-d12	13.84

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	94
2		Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	93
3		Pentafluoropropionic acid, tetra...	360	C17H29F5O2	006222-06-6	90
4		Pentafluoropropionic acid, penta...	374	C18H31F5O2	959092-08-1	90
5		n-Tetracosanol-1	354	C24H50O	000506-51-4	90



Data Path : Z:\HPCHEM1\BNA F\DATA\BF071216\
Data File : BF088940.D
Acq On : 13 Jul 2016 1:24
Operator : UM/SJ
Sample : H3935-27
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
C3-01

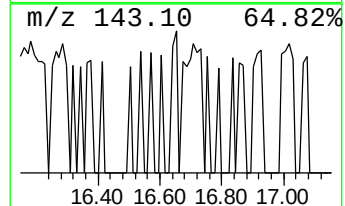
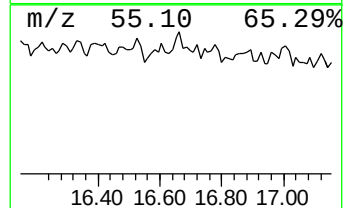
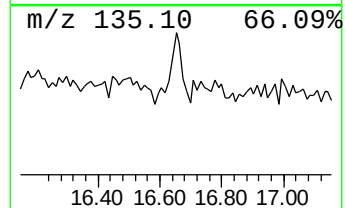
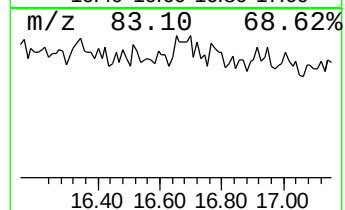
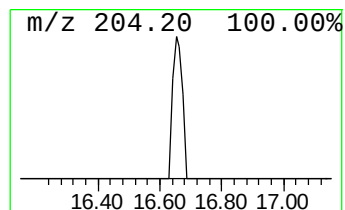
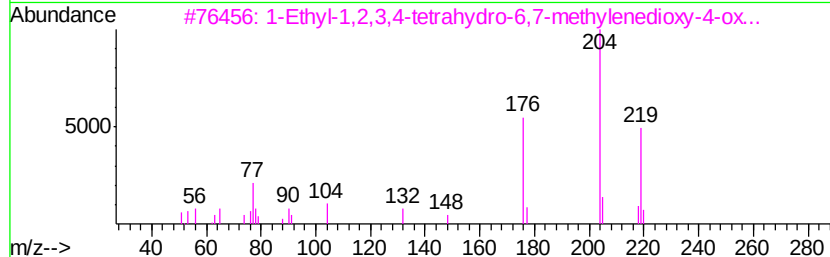
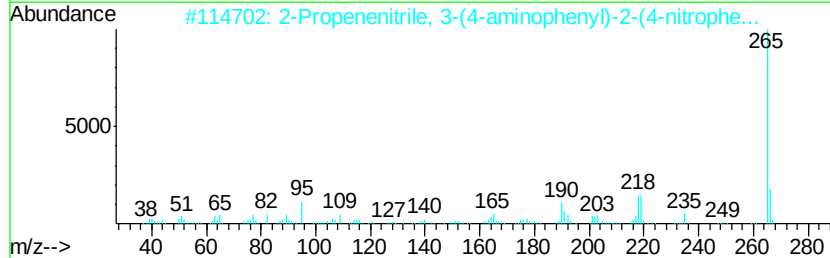
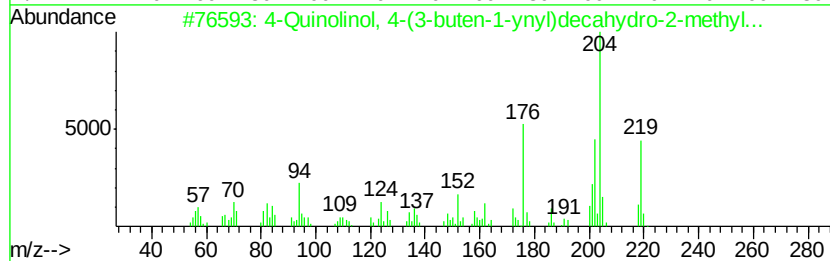
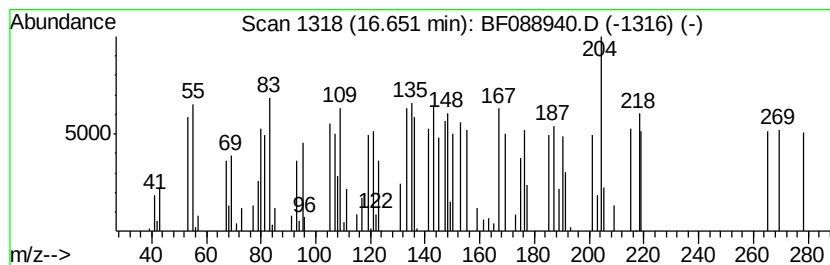
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 16 unknown16.65 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.65	2.42 ng	40633	Perylene-d12	15.21

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-Quinolinol, 4-(3-buten-1-ynyl)...	219	C14H21NO	038423-21-1	25
2			2-Propenenitrile, 3-(4-aminophen...	265	C15H11N3O2	013491-58-2	14
3			1-Ethyl-1,2,3,4-tetrahydro-6,7-m...	219	C12H13NO3	132767-88-5	10
4			9b-Boraphenalene, dodecahydro-	176	C12H21B	016664-33-8	10
5			2-Oxazoline, 4,5-diethyl-2-(p-ni...	248	C13H16N2O3	025943-08-2	10



Data Path : Z:\HPCHEM1\BNA F\DATA\BF071216\
Data File : BF088940.D
Acq On : 13 Jul 2016 1:24
Operator : UM/SJ
Sample : H3935-27
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
C3-01

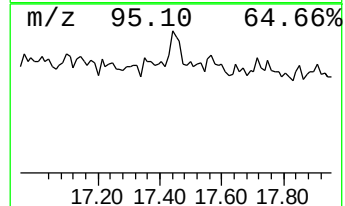
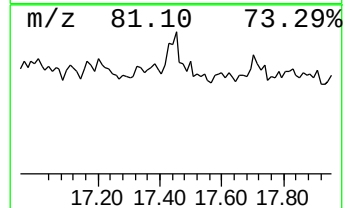
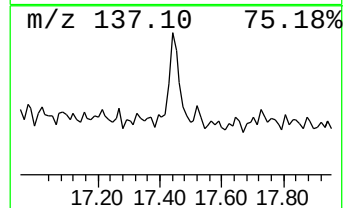
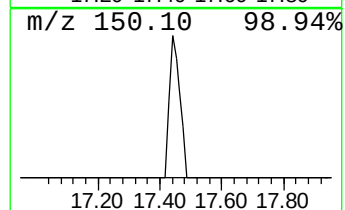
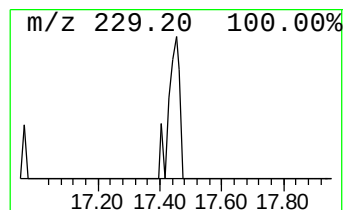
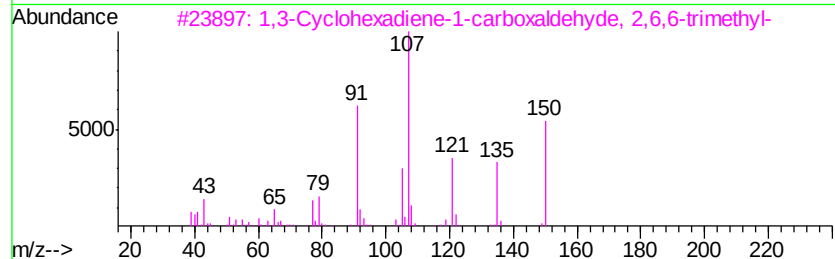
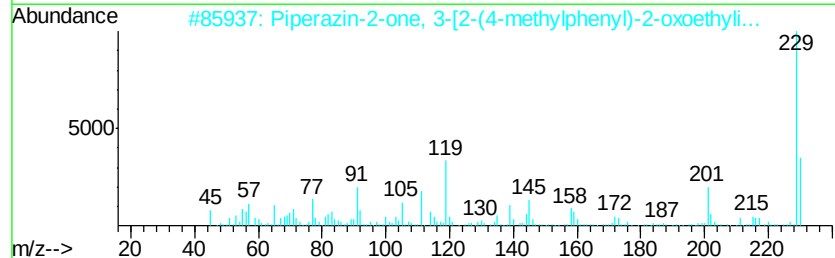
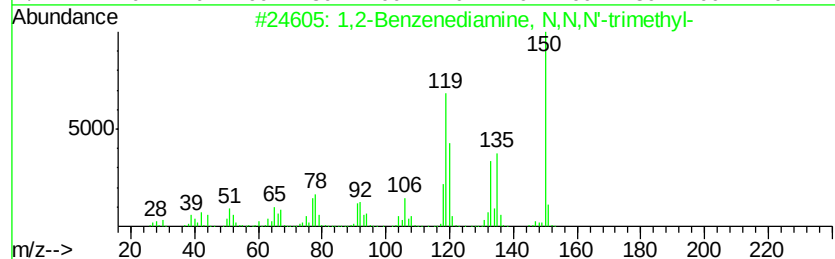
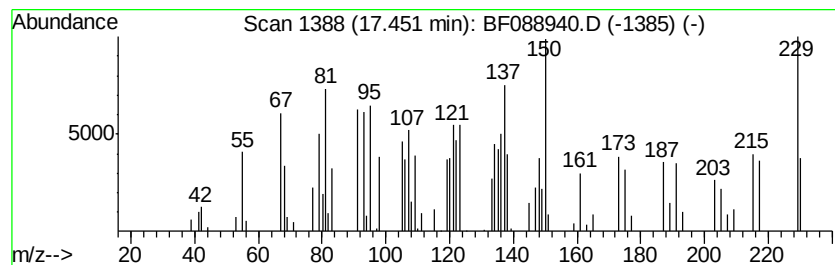
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF063016.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 17 unknown17.45 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.45	2.85 ng	47941	Perylene-d12	15.21

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,2-Benzenediamine, N,N,N'-trime...	150	C9H14N2	002427-03-4	11
2			Piperazin-2-one, 3-[2-(4-methylp...	230	C13H14N2O2	063656-21-3	10
3			1,3-Cyclohexadiene-1-carboxaldeh...	150	C10H14O	000116-26-7	10
4			3-Methoxy-4-methylaniline	137	C8H11NO	016452-01-0	10
5			2-Cyclopenten-1-one, 2-(2-buteny...	150	C10H14O	017190-71-5	10



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF071216\
 Data File : BF088940.D
 Acq On : 13 Jul 2016 1:24
 Operator : UM/SJ
 Sample : H3935-27
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 C3-01

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF063016.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Propane, 1,1-dime...	1.88	4.6	ng	92681	1	6.71	400693	20.0
unknown2.86	2.86	2.5	ng	51129	1	6.71	400693	20.0
2-Pentanone, 4-hy...	4.86	11.7	ng	233569	1	6.71	400693	20.0
unknown6.48	6.48	100.9	ng	2020720	1	6.71	400693	20.0
Tridecane	10.65	3.2	ng	64011	4	11.21	404192	20.0
1-Tetradecene	10.97	5.6	ng	113235	4	11.21	404192	20.0
Dodecane	11.10	3.1	ng	63004	4	11.21	404192	20.0
1-Hexadecanol	11.45	2.1	ng	42927	4	11.21	404192	20.0
unknown11.76	11.76	2.0	ng	41334	4	11.21	404192	20.0
unknown11.82	11.82	3.0	ng	59855	4	11.21	404192	20.0
1-Heneicosanol	13.70	7.2	ng	148904	5	13.84	413115	20.0
Heptadecyl heptaf...	14.33	5.4	ng	110892	5	13.84	413115	20.0
unknown16.65	16.65	2.4	ng	40633	6	15.21	336438	20.0
unknown17.45	17.45	2.9	ng	47941	6	15.21	336438	20.0