

Data Path : Z:\HPCHEM1\BNA F\DATA\BF052215\
 Data File : BF079468.D
 Acq On : 23 May 2015 10:47
 Operator : TP/IZ
 Sample : G2371-01
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SW-021

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	1.324	8	12	15	rVB	370557	699110	16.94%	2.809%
2	1.484	24	26	29	rVB	246648	281034	6.81%	1.129%
3	1.610	35	37	45	rVB	220273	432293	10.47%	1.737%
4	1.713	45	46	83	rVB	2183950	4128047	100.00%	16.585%
5	4.787	313	315	322	rVB	103558	146084	3.54%	0.587%
6	5.347	361	364	376	rBV	1418988	1664042	40.31%	6.685%
7	6.650	476	478	486	rVB	1739147	1797387	43.54%	7.221%
8	6.765	486	488	491	rBV	1332588	1786082	43.27%	7.176%
9	7.050	511	513	516	rBV	510794	462181	11.20%	1.857%
10	7.233	527	529	532	rBV	1346607	1376917	33.36%	5.532%
11	7.748	572	574	577	rBV	821776	720063	17.44%	2.893%
12	8.376	616	629	639	rBV	49825	182847	4.43%	0.735%
13	8.628	649	651	654	rBV	720189	640028	15.50%	2.571%
14	9.874	758	760	765	rBV	205936	217445	5.27%	0.874%
15	9.976	767	769	771	rBV	2482323	2199435	53.28%	8.836%
16	10.491	811	814	816	rBV	119763	151180	3.66%	0.607%
17	10.799	836	841	843	rBV2	632692	798346	19.34%	3.207%
18	11.782	924	927	929	rBV2	518686	763935	18.51%	3.069%
19	12.445	982	985	987	rBV	63279	85817	2.08%	0.345%
20	12.628	999	1001	1003	rBV	613736	721689	17.48%	2.899%
21	13.371	1063	1066	1069	rBV	150753	235993	5.72%	0.948%
22	13.908	1110	1113	1118	rBV	59092	108745	2.63%	0.437%
23	14.137	1128	1133	1135	rBV	68060	100632	2.44%	0.404%
24	14.263	1140	1144	1146	rBV4	21478	53027	1.28%	0.213%
25	14.411	1155	1157	1159	rBV	60258	63730	1.54%	0.256%
26	14.628	1173	1176	1179	rVB	2283133	2309044	55.94%	9.277%
27	14.720	1179	1184	1188	rBV3	49792	94600	2.29%	0.380%
28	15.200	1223	1226	1229	rBV	24580	57746	1.40%	0.232%
29	15.337	1236	1238	1240	rBV2	27503	56607	1.37%	0.227%
30	15.806	1276	1279	1282	rBV	212822	346615	8.40%	1.393%
31	15.874	1282	1285	1286	rBV2	39048	69185	1.68%	0.278%
32	15.908	1286	1288	1291	rVV	540895	786557	19.05%	3.160%
33	16.080	1300	1303	1305	rBV	116511	138719	3.36%	0.557%
34	16.617	1348	1350	1353	rVB2	38833	58390	1.41%	0.235%

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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

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Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	16.686	1353	1356	1357	rBV3	28169	50598	1.23%	0.203%
36	16.743	1358	1361	1365	rBV	47535	156976	3.80%	0.631%
37	16.880	1371	1373	1377	rVB4	36155	55463	1.34%	0.223%
38	16.960	1377	1380	1382	rBV2	51279	89071	2.16%	0.358%
39	17.646	1437	1440	1443	rBV	663422	805247	19.51%	3.235%

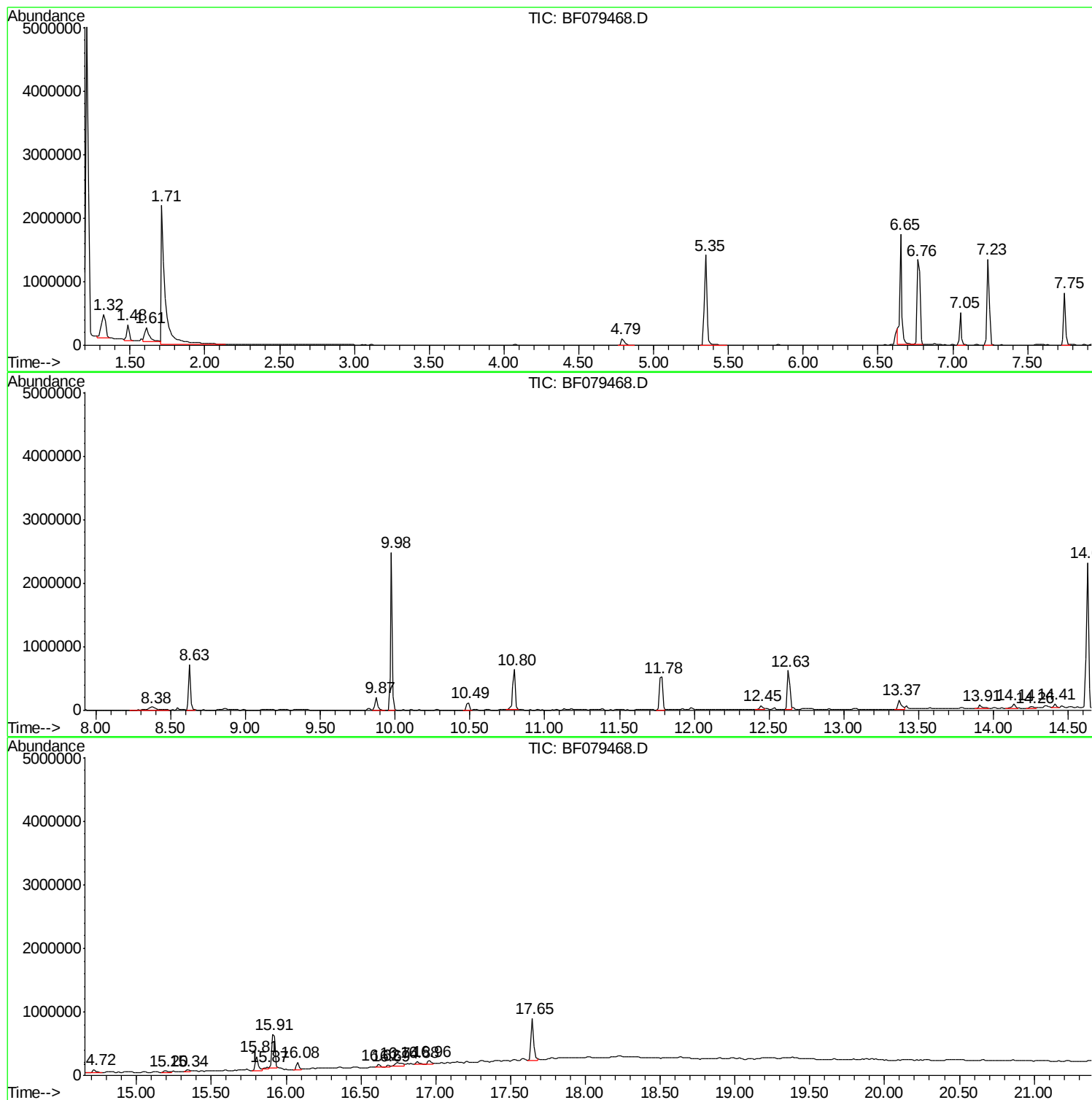
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF043015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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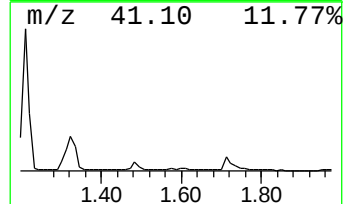
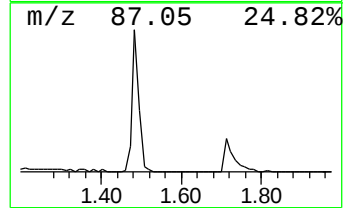
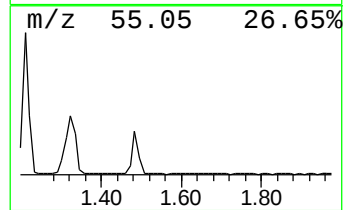
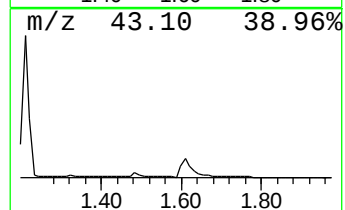
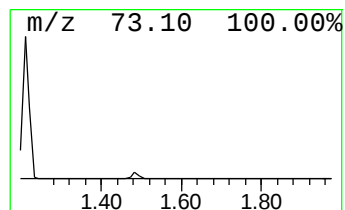
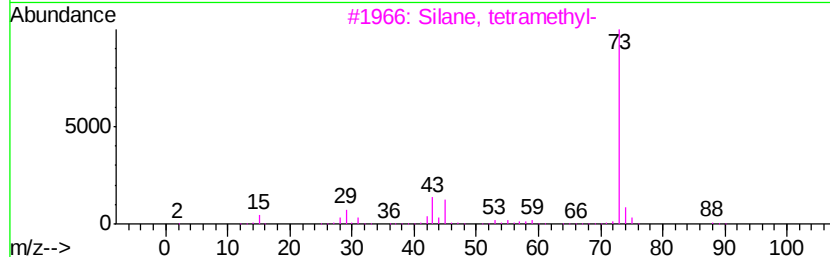
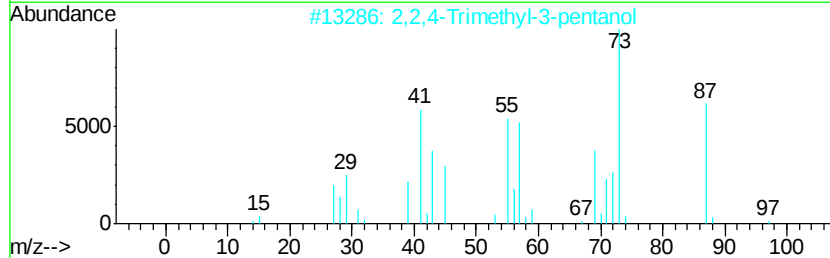
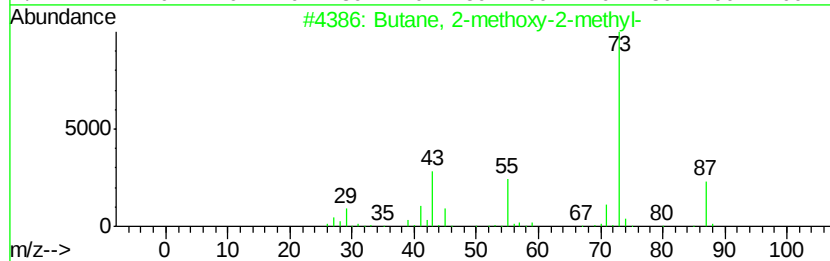
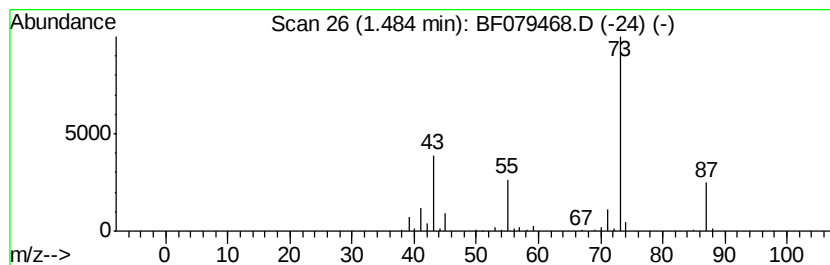
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Butane, 2-methoxy-2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.48	12.16 ng	281034	1,4-Dichlorobenzene-d4	7.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	72
2		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
3		Silane, tetramethyl-	88	C4H12Si	000075-76-3	35
4		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12
5		N-Ethylformamide	73	C3H7NO	000627-45-2	9



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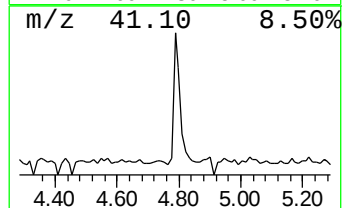
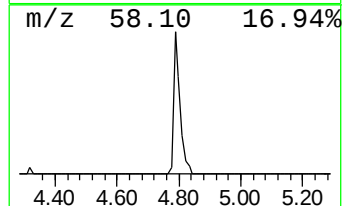
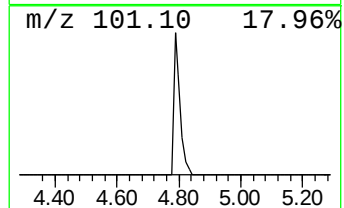
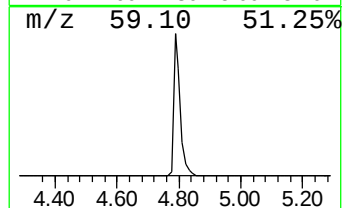
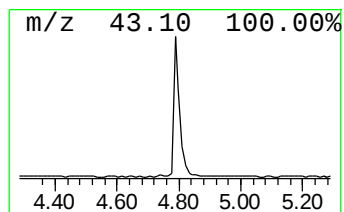
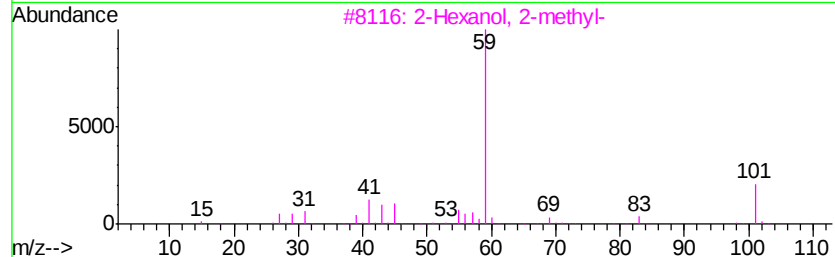
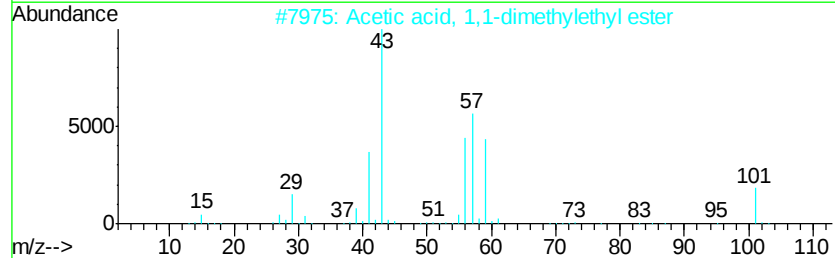
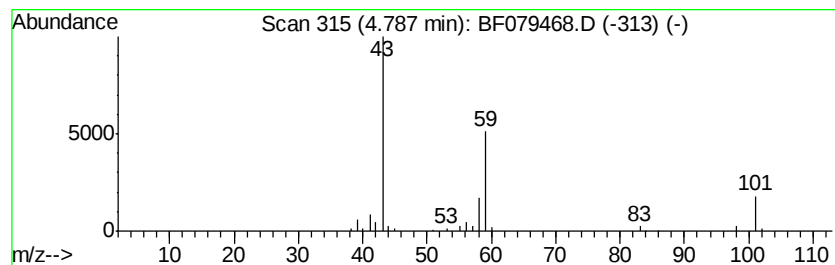
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.79	6.32 ng	146084	1,4-Dichlorobenzene-d4	7.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	38
3		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	33
4		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	10
5		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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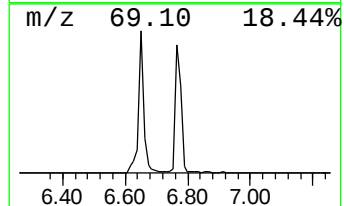
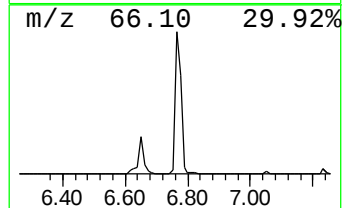
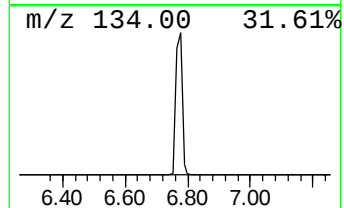
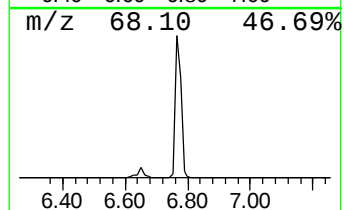
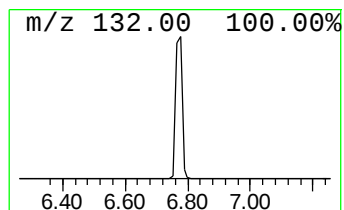
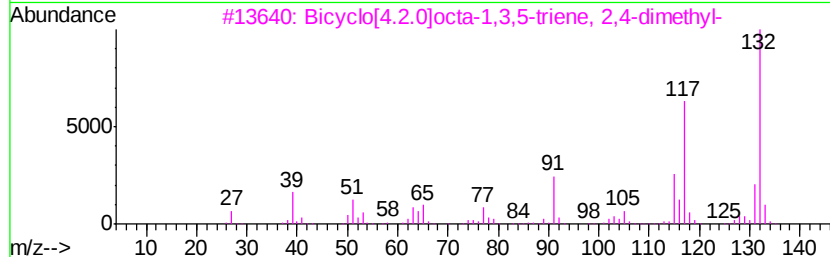
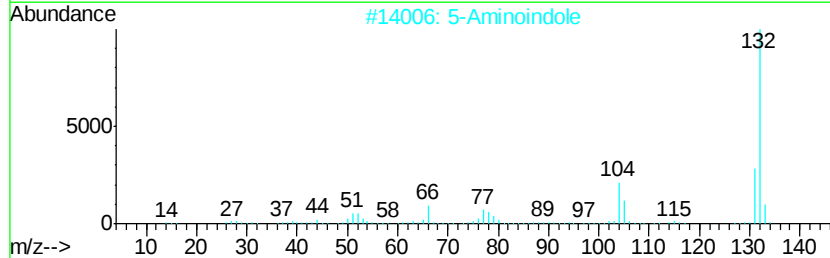
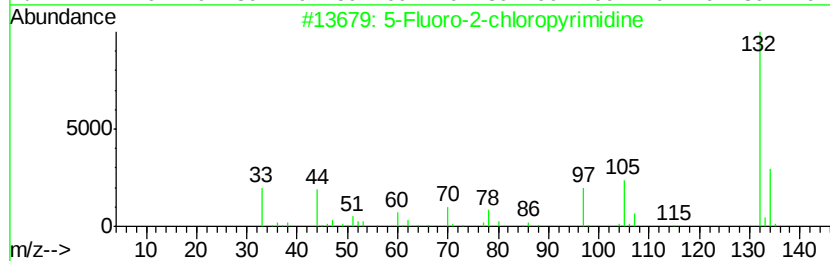
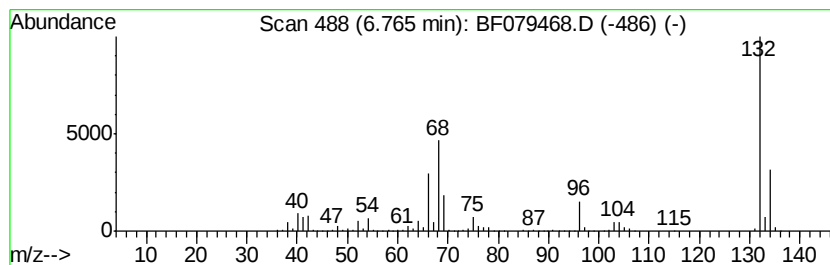
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 unknown6.76 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.76	77.29 ng	1786080	1,4-Dichlorobenzene-d4	7.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
2		5-Aminoindole	132	C8H8N2	005192-03-0	12
3		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
4		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9
5		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9



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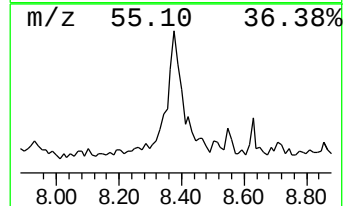
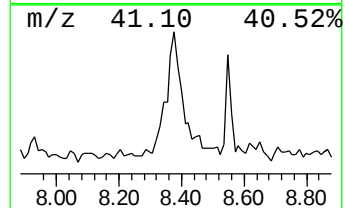
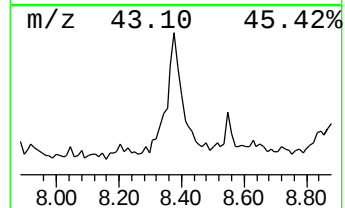
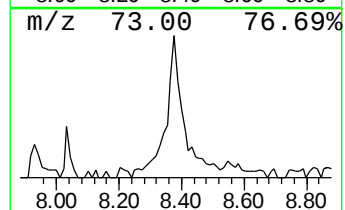
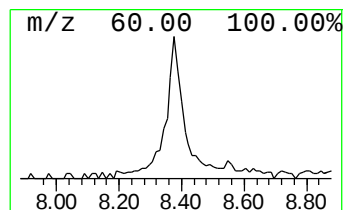
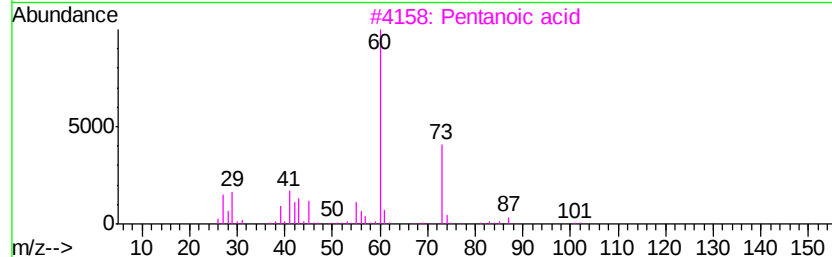
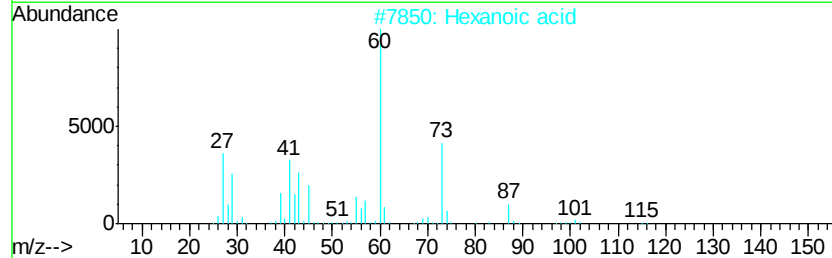
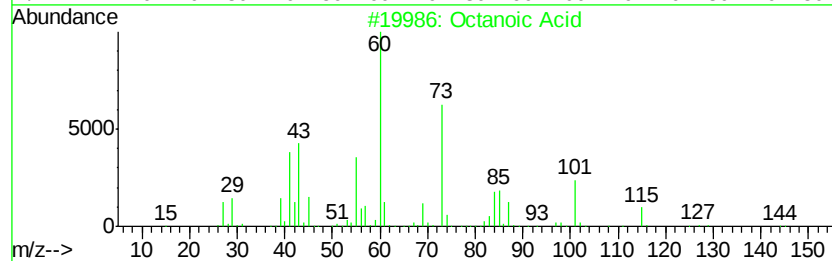
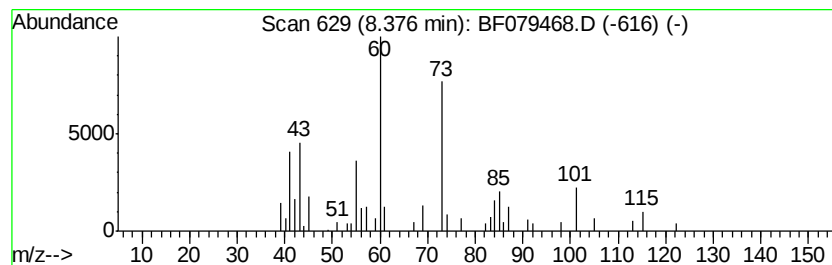
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Octanoic Acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.38	5.71 ng	182847	Naphthalene-d8	8.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octanoic Acid	144	C8H16O2	000124-07-2	91
2		Hexanoic acid	116	C6H12O2	000142-62-1	50
3		Pentanoic acid	102	C5H10O2	000109-52-4	40
4		Heptanoic acid	130	C7H14O2	000111-14-8	37
5		Propanedioic acid, propyl-	146	C6H10O4	000616-62-6	37



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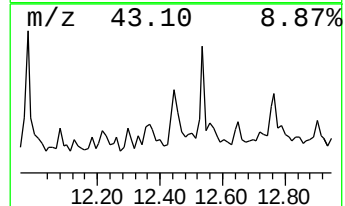
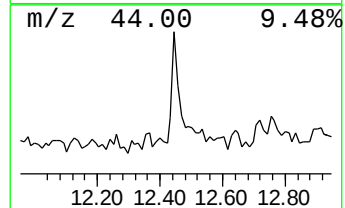
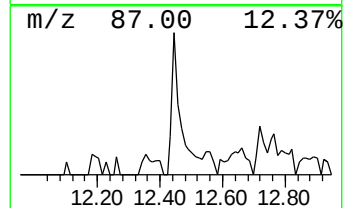
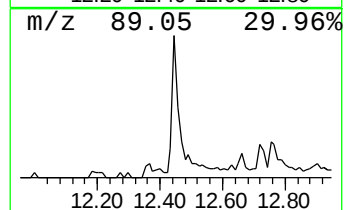
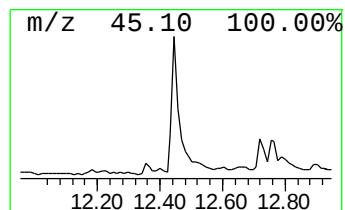
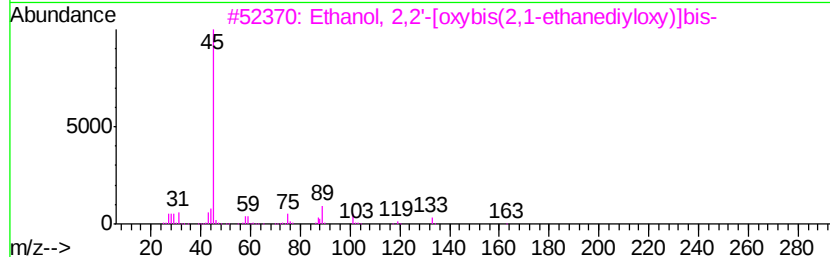
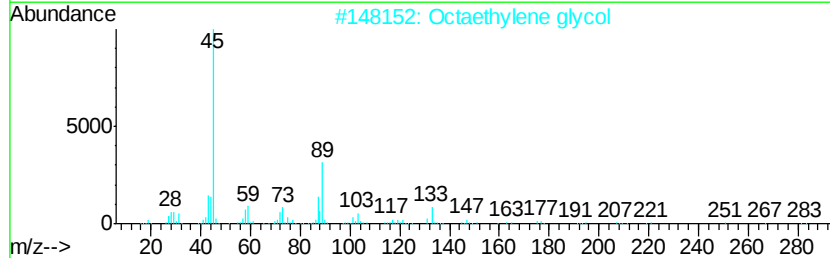
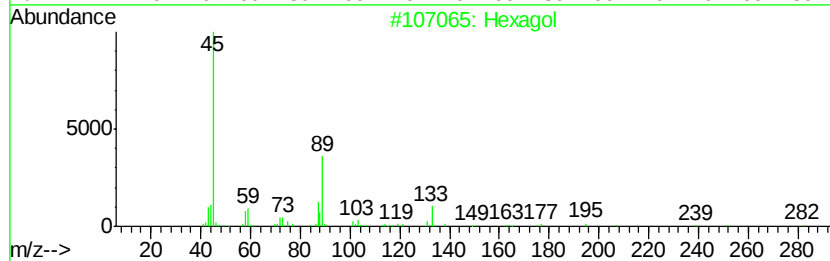
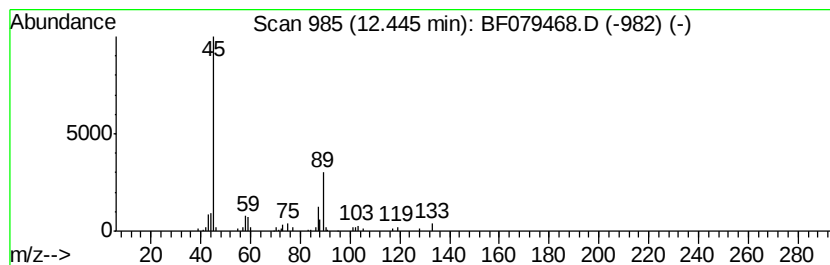
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Hexagol Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.45	2.38 ng	85817	Phenanthrene-d10	12.63
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Hexagol		282 C12H26O7	002615-15-8 83
2	Octaethylene glycol		370 C16H34O9	1000289-34-2 74
3	Ethanol, 2,2'-[oxybis(2,1-ethane...		194 C8H18O5	000112-60-7 72
4	Triethylene glycol		150 C6H14O4	000112-27-6 64
5	2,5,8,11,14-Pentaoxahexadecan-16-ol		252 C11H24O6	023778-52-1 64



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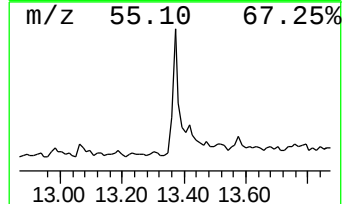
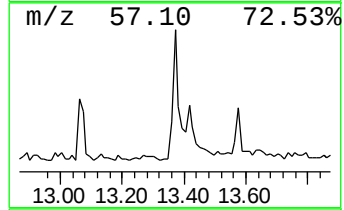
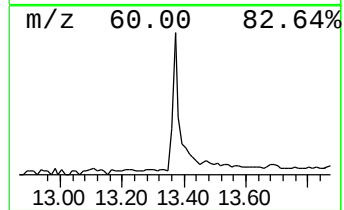
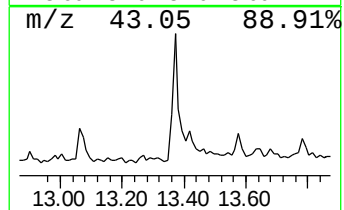
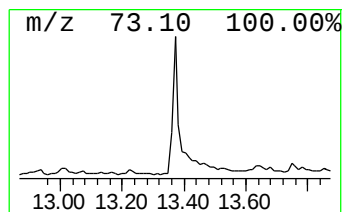
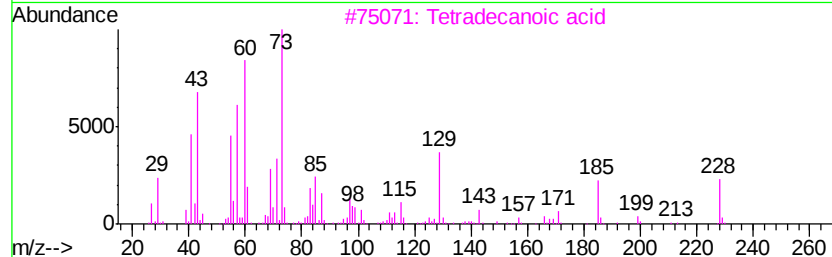
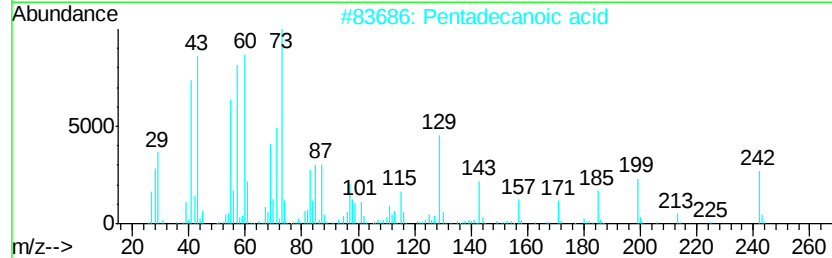
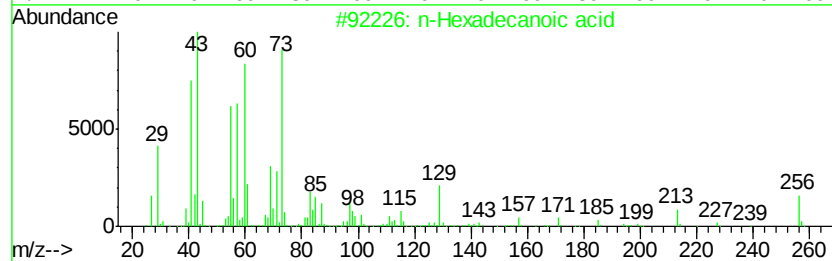
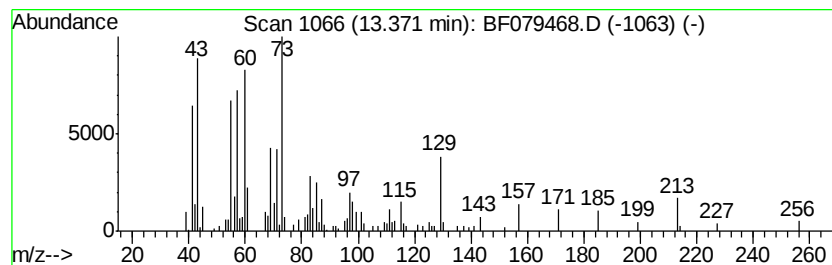
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 n-Hexadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.37	6.54 ng	235993	Phenanthrene-d10	12.63

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	91
2			Pentadecanoic acid	242	C15H30O2	001002-84-2	87
3			Tetradecanoic acid	228	C14H28O2	000544-63-8	83
4			Tridecanoic acid	214	C13H26O2	000638-53-9	76
5			n-Decanoic acid	172	C10H20O2	000334-48-5	64



Data Path : Z:\HPCHEM1\BNA F\DATA\BF052215\
Data File : BF079468.D
Acq On : 23 May 2015 10:47
Operator : TP/IZ
Sample : G2371-01
Misc :
ALS Vial : 36 Sample Multiplier: 1

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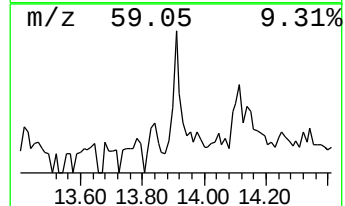
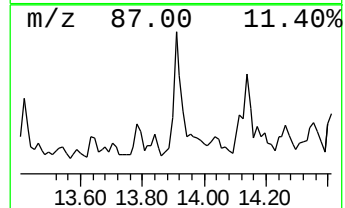
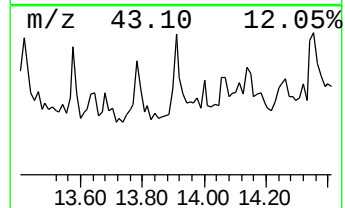
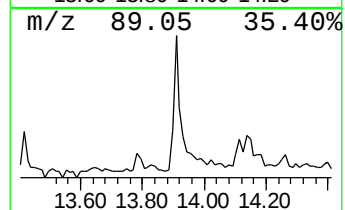
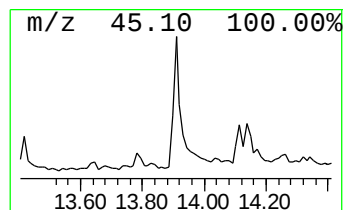
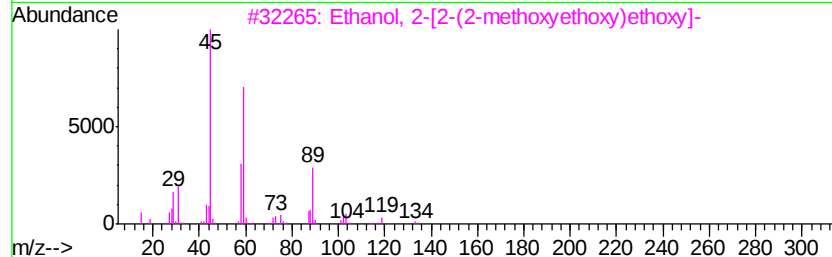
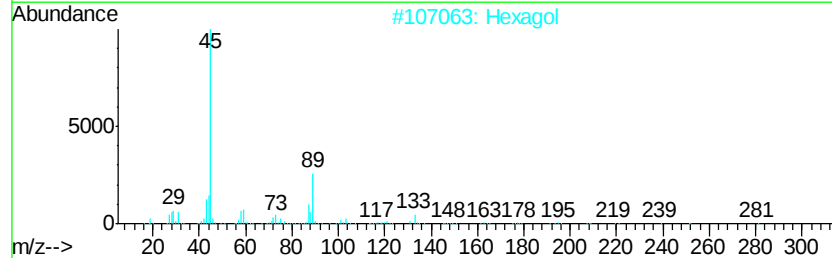
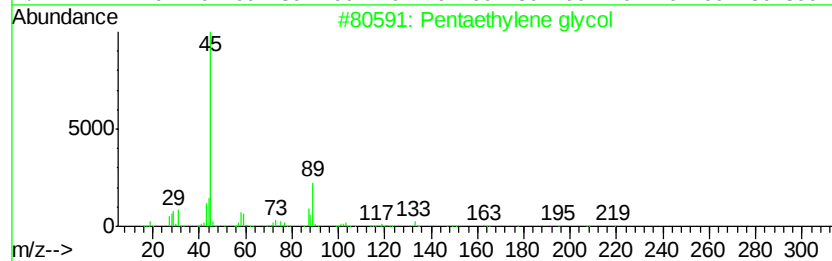
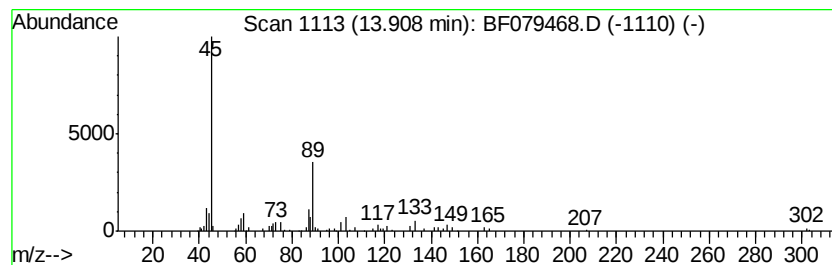
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Pentaethylene glycol Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.91	3.01 ng	108745	Phenanthrene-d10	12.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentaethylene glycol	238	C10H22O6	004792-15-8	83
2		Hexagol	282	C12H26O7	002615-15-8	83
3		Ethanol, 2-[2-(2-methoxyethoxy)ethoxy]ethoxy]	164	C7H16O4	000112-35-6	72
4		Ethanol, 2,2'-[oxybis(2,1-ethane)]	194	C8H18O5	000112-60-7	72
5		Acetaldehyde, tetramer	176	C8H16O4	000108-62-3	59



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF052215\
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Misc :
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ClientSampleId :
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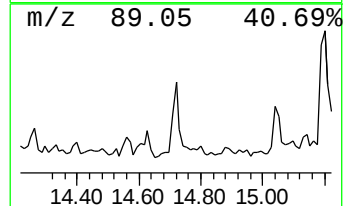
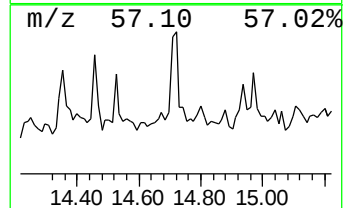
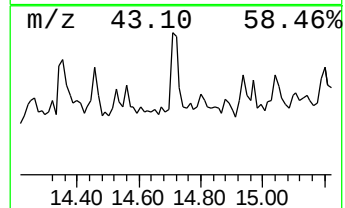
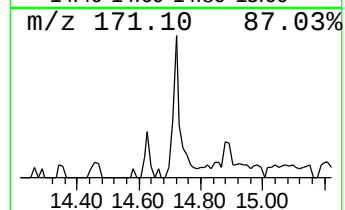
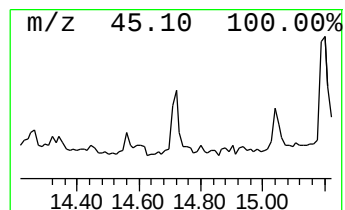
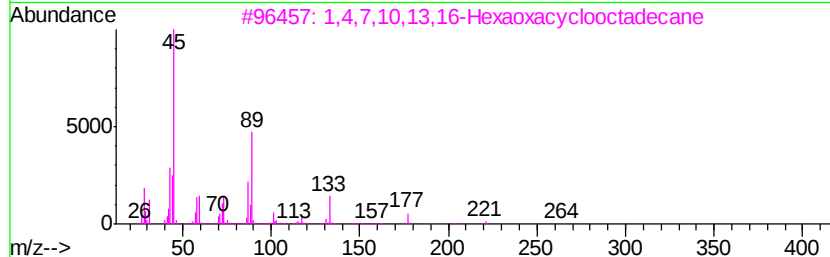
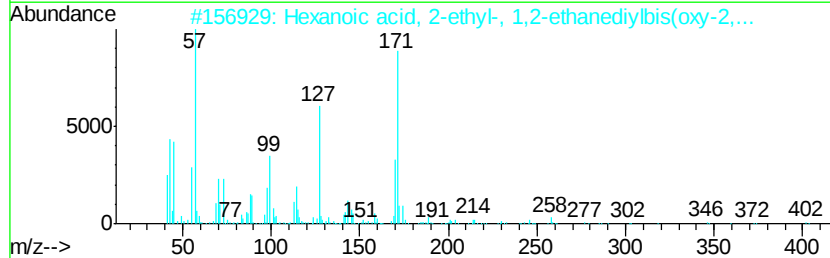
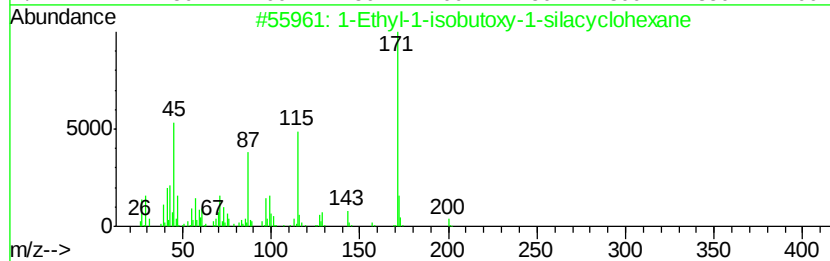
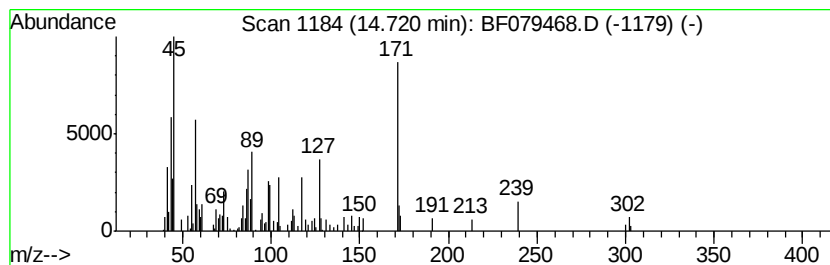
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 unknown14.72 Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.72	2.41 ng	94600	Chrysene-d12	15.92

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Ethyl-1-isobutoxy-1-silacycloh...	200	C11H24OSi	1000279-31-1	47
2		Hexanoic acid, 2-ethyl-, 1,2-eth...	402	C22H42O6	000094-28-0	37
3		1,4,7,10,13,16-Hexaoxacyclooctad...	264	C12H24O6	017455-13-9	27
4		Dimethyl-dioctyl-silane	284	C18H40Si	003429-75-2	27
5		Hexagol	282	C12H26O7	002615-15-8	27



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Sample : G2371-01
Misc :
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ClientSampleId :
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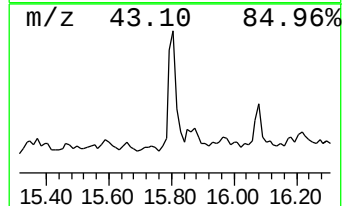
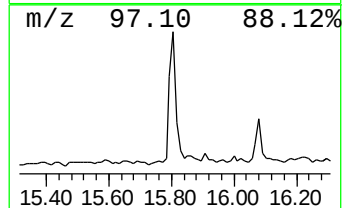
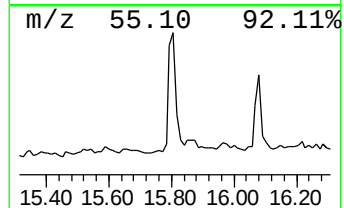
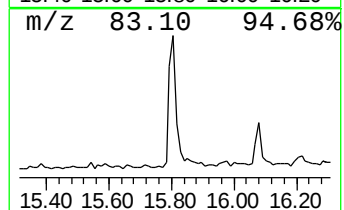
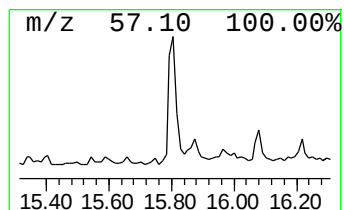
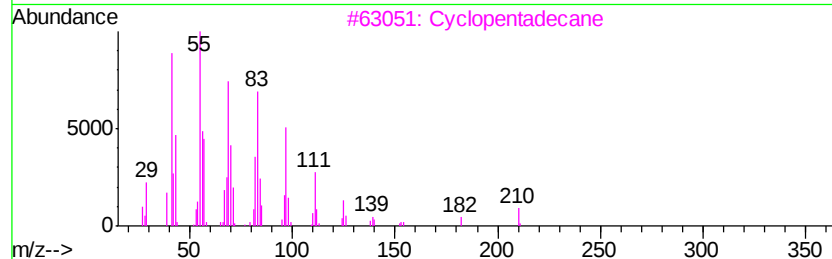
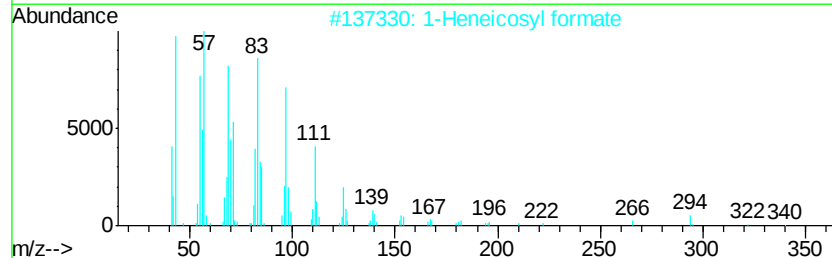
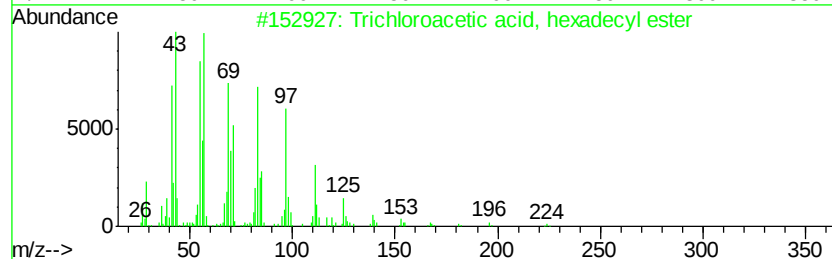
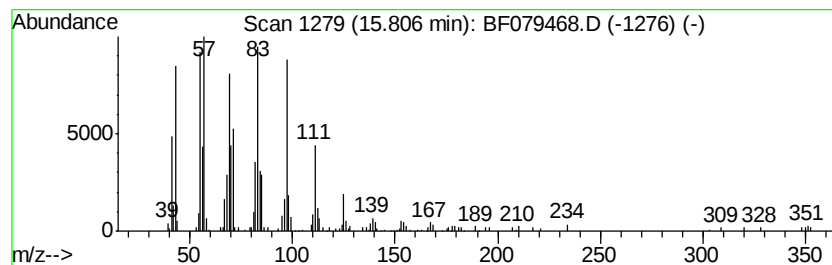
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 15 Trichloroacetic acid, hexad... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.81	8.81 ng	346615	Chrysene-d12	15.92

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	94
2		1-Heneicosyl formate	340	C22H44O2	077899-03-7	93
3		Cyclopentadecane	210	C15H30	000295-48-7	93
4		1-Nonadecene	266	C19H38	018435-45-5	91
5		9-Eicosene, (E)-	280	C20H40	074685-29-3	91



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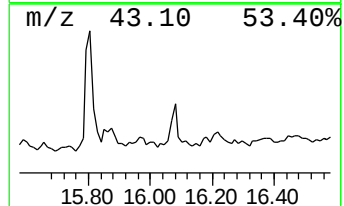
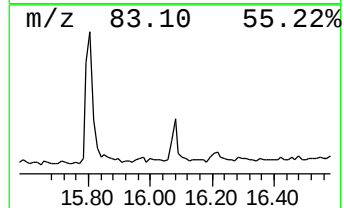
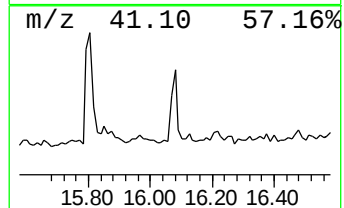
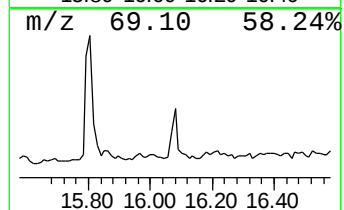
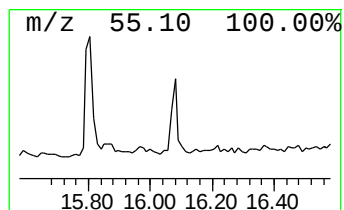
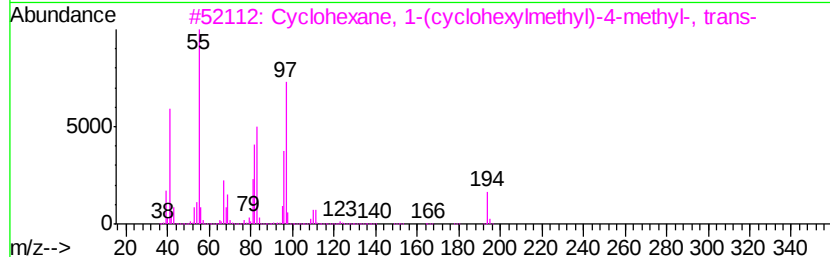
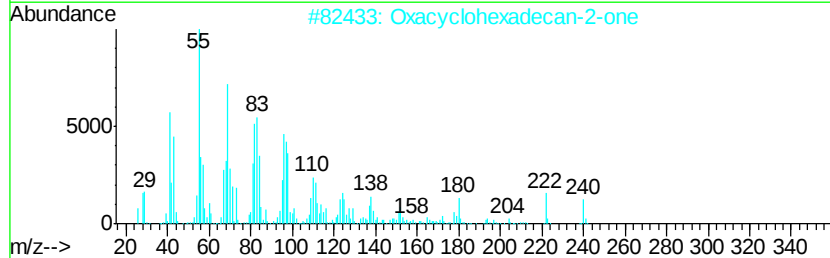
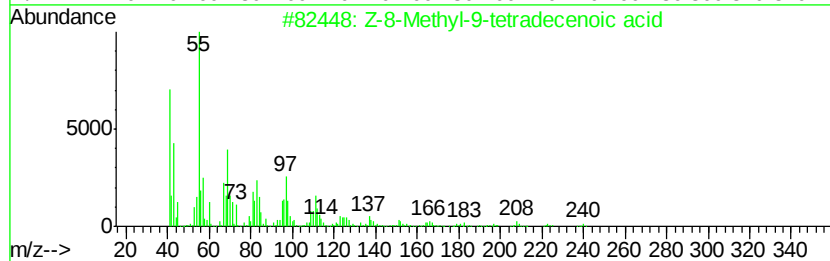
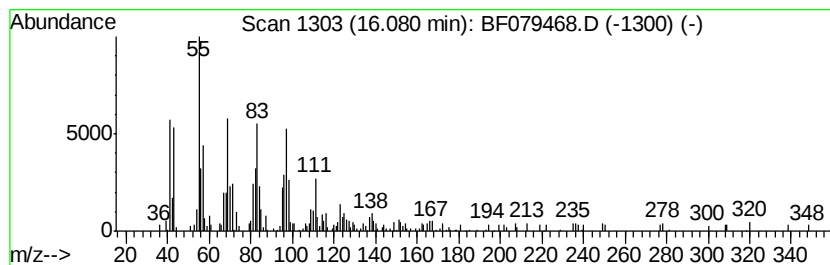
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 16 Z-8-Methyl-9-tetradecenoic ... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.08	3.53 ng	138719	Chrysene-d12	15.92

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Z-8-Methyl-9-tetradecenoic acid	240	C15H28O2	1000130-84-5	81
2			Oxacyclohexadecan-2-one	240	C15H28O2	000106-02-5	60
3			Cyclohexane, 1-(cyclohexylmethyl)-	194	C14H26	054823-98-2	60
4			Oxacyclotetradecane-2,11-dione, ...	240	C14H24O3	074685-36-2	53
5			Cyclopentadecanone, 2-hydroxy-	240	C15H28O2	004727-18-8	49



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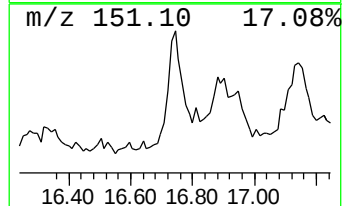
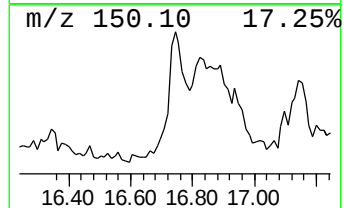
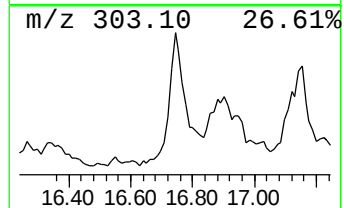
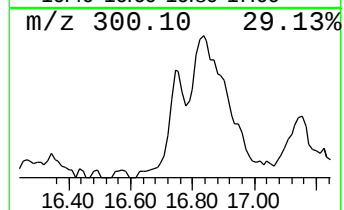
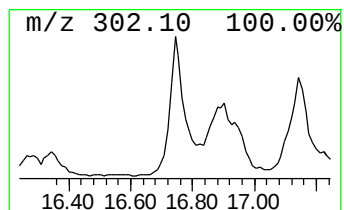
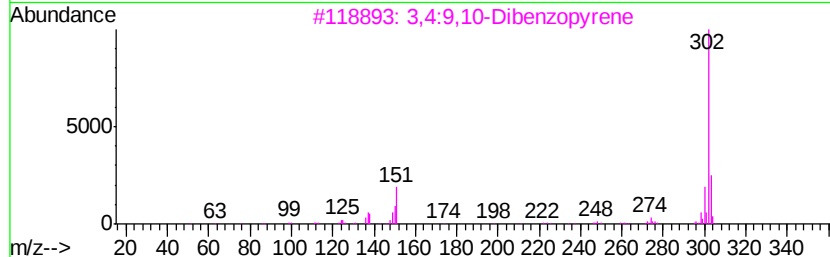
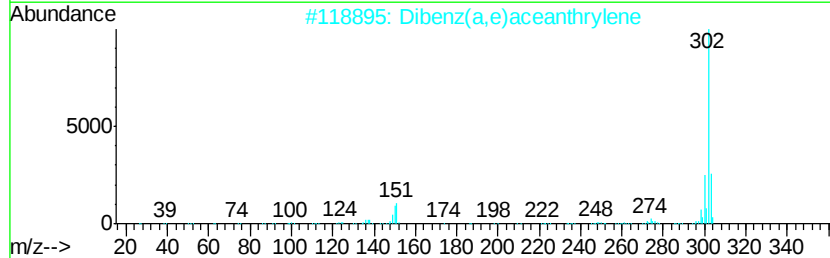
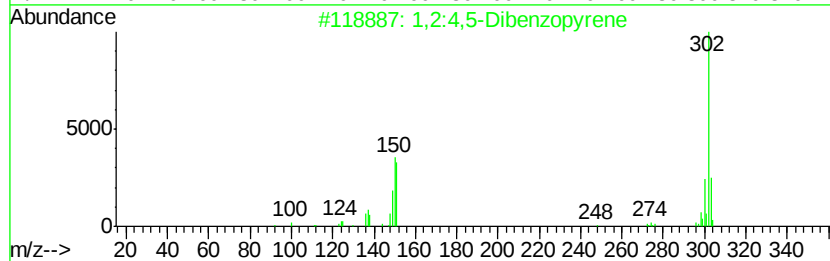
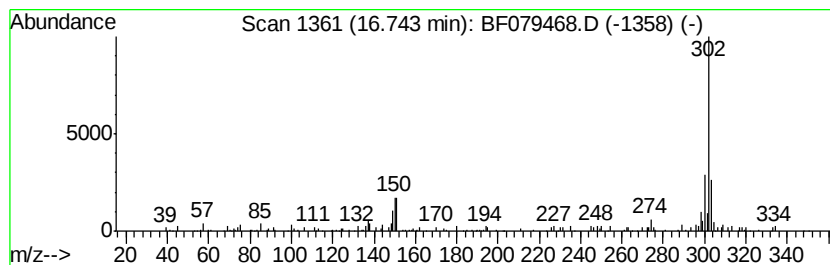
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 1,2:4,5-Dibenzopyrene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.74	3.99 ng	156976	Chrysene-d12	15.92

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,2:4,5-Dibenzopyrene	302	C24H14	000192-65-4	95
2			Dibenz(a,e)aceanthrylene	302	C24H14	005385-75-1	93
3			3,4:9,10-Dibenzopyrene	302	C24H14	000189-55-9	93
4			Indeno[1,2,3-f]naphthacene	302	C24H14	000203-11-2	90
5			3,4:8,9-Dibenzopyrene	302	C24H14	000189-64-0	90



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ALS Vial : 36 Sample Multiplier: 1

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-methoxy...	1.48	12.2	ng	281034	1	7.05	462181	20.0
2-Pentanone, 4-hy...	4.79	6.3	ng	146084	1	7.05	462181	20.0
unknown6.76	6.76	77.3	ng	1786080	1	7.05	462181	20.0
Octanoic Acid	8.38	5.7	ng	182847	2	8.63	640028	20.0
Hexagol	12.45	2.4	ng	85817	4	12.63	721689	20.0
n-Hexadecanoic acid	13.37	6.5	ng	235993	4	12.63	721689	20.0
Pentaethylene glycol	13.91	3.0	ng	108745	4	12.63	721689	20.0
unknown14.72	14.72	2.4	ng	94600	5	15.92	786557	20.0
Trichloroacetic a...	15.81	8.8	ng	346615	5	15.92	786557	20.0
Z-8-Methyl-9-tetr...	16.08	3.5	ng	138719	5	15.92	786557	20.0
1,2:4,5-Dibenzopy...	16.74	4.0	ng	156976	5	15.92	786557	20.0