

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF021716\
 Data File : BF085095.D
 Acq On : 17 Feb 2016 19:20
 Operator : SJ/IZ
 Sample : H1443-01
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EP004

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-BF021716.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	1.873	22	25	28	rVB	5206446	6195475	100.00%	11.315%
2	2.010	34	37	46	rVB2	1473492	2580510	41.65%	4.713%
3	2.124	46	47	52	rVB	52828	79055	1.28%	0.144%
4	2.204	52	54	60	rVB	108818	163182	2.63%	0.298%
5	2.330	63	65	79	rVB3	32145	88431	1.43%	0.162%
6	5.187	313	315	318	rVB	488384	652841	10.54%	1.192%
7	5.599	348	351	353	rBV	2622672	3164631	51.08%	5.780%
8	6.605	436	439	442	rBV	2607884	3443966	55.59%	6.290%
9	6.753	450	452	455	rBV	2633447	3354869	54.15%	6.127%
10	6.993	471	473	475	rBV	1464667	1182890	19.09%	2.160%
11	7.153	484	487	489	rVB	2796615	2684769	43.33%	4.903%
12	7.553	519	522	524	rBV	2570390	2280703	36.81%	4.165%
13	8.273	583	585	588	rBV	1187972	1519846	24.53%	2.776%
14	9.348	677	679	682	rVB	3555515	4877672	78.73%	8.908%
15	9.748	712	714	715	rBV	668768	734599	11.86%	1.342%
16	9.965	730	733	735	rVB	142387	138179	2.23%	0.252%
17	10.034	736	739	741	rVB	2104551	1773469	28.63%	3.239%
18	10.456	775	776	778	rVB	209666	220257	3.56%	0.402%
19	10.685	793	796	799	rBV	80360	153570	2.48%	0.280%
20	10.822	805	808	810	rBV	3805841	3954564	63.83%	7.222%
21	10.936	816	818	822	rBV	239988	320586	5.17%	0.585%
22	11.382	854	857	858	rBV	133078	125753	2.03%	0.230%
23	11.519	866	869	870	rBV	1870729	1855648	29.95%	3.389%
24	11.542	870	871	872	rVV	306345	219449	3.54%	0.401%
25	11.588	872	875	877	rVB	73297	93236	1.50%	0.170%
26	12.045	910	915	917	rBV2	250908	468196	7.56%	0.855%
27	12.125	920	922	927	rVB2	94836	157696	2.55%	0.288%
28	12.331	934	940	941	rBV2	59583	176290	2.85%	0.322%
29	12.457	943	951	952	rBV3	70586	286278	4.62%	0.523%
30	12.651	966	968	970	rVB3	43605	62145	1.00%	0.113%
31	12.731	973	975	977	rVB	212935	289347	4.67%	0.528%
32	12.822	981	983	987	rBV2	120051	160566	2.59%	0.293%
33	12.948	992	994	998	rVB	207500	336100	5.42%	0.614%
34	13.097	1005	1007	1010	rVB	4053788	5544288	89.49%	10.126%

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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

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Peak separation: 5

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35	13.177	1010	1014	1015	rBV	38900	68655	1.11%	0.125%
36	13.280	1020	1023	1027	rVV2	93024	254714	4.11%	0.465%
37	13.382	1030	1032	1036	rVB2	43143	67811	1.09%	0.124%
38	13.988	1082	1085	1087	rBV	866666	790796	12.76%	1.444%
39	14.148	1096	1099	1104	rVB	1539087	1640350	26.48%	2.996%
40	14.617	1138	1140	1143	rBV3	128688	246649	3.98%	0.450%
41	14.937	1165	1168	1172	rVV4	60669	152896	2.47%	0.279%
42	15.188	1188	1190	1195	rVB	74218	147252	2.38%	0.269%
43	15.497	1215	1217	1220	rVB	55332	66013	1.07%	0.121%
44	15.554	1220	1222	1225	rBV	57634	89705	1.45%	0.164%
45	15.623	1225	1228	1231	rBV	818537	1125433	18.17%	2.055%
46	15.988	1258	1260	1269	rVB2	79051	214003	3.45%	0.391%
47	16.148	1272	1274	1283	rVB2	110282	219493	3.54%	0.401%
48	16.377	1290	1294	1299	rVB3	69066	197777	3.19%	0.361%
49	17.097	1354	1357	1361	rVB2	32440	69396	1.12%	0.127%
50	17.314	1373	1376	1380	rVB3	28736	64581	1.04%	0.118%

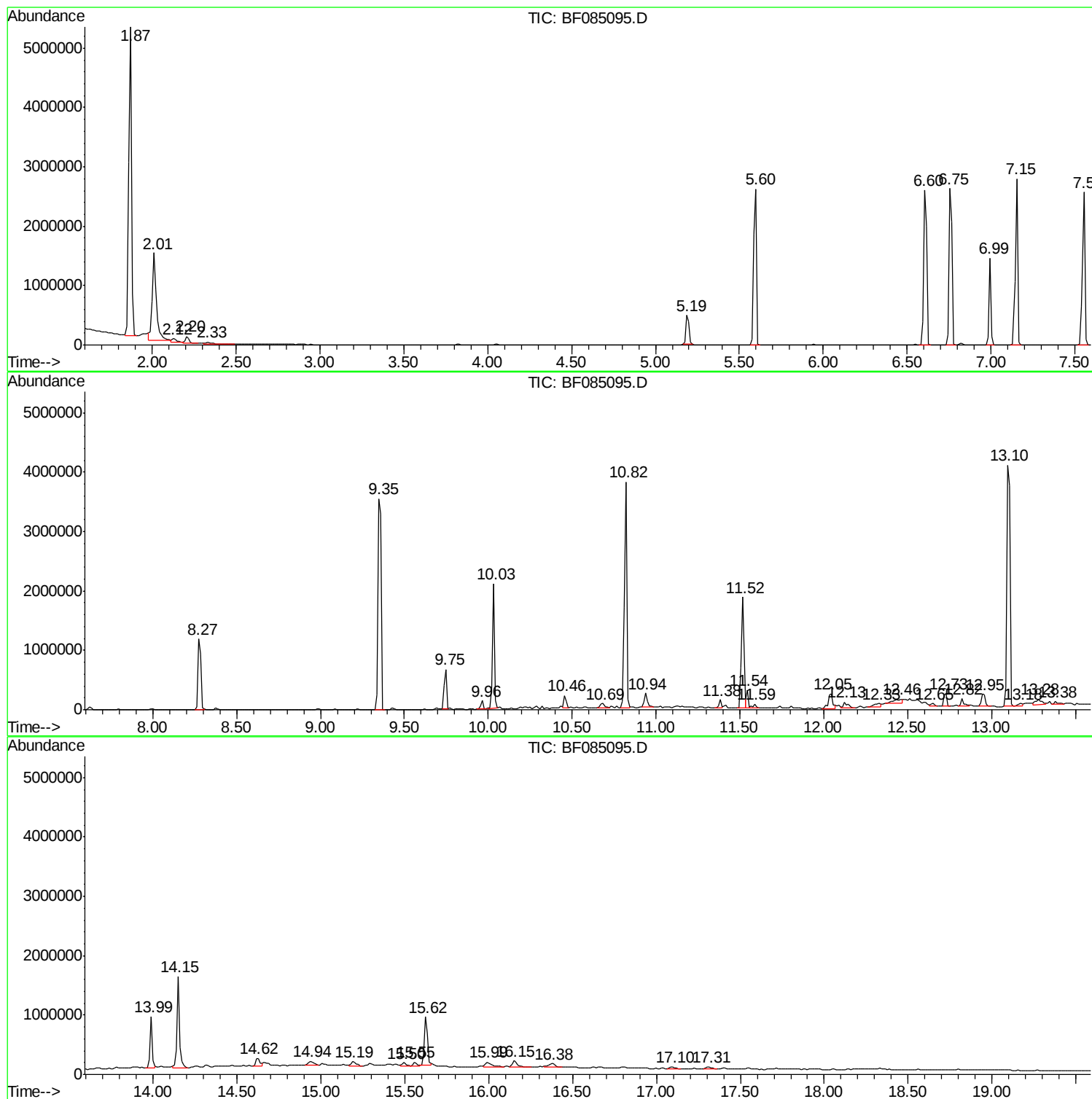
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ClientSampled :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF021716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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Sample : H1443-01
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Instrument :
BNA_F
ClientSampleId :
EP004

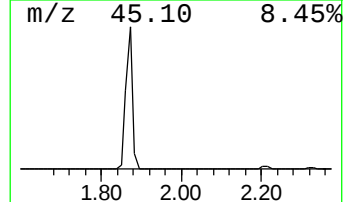
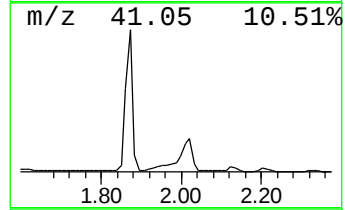
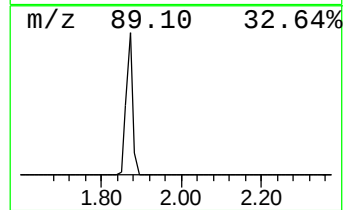
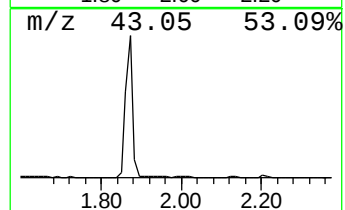
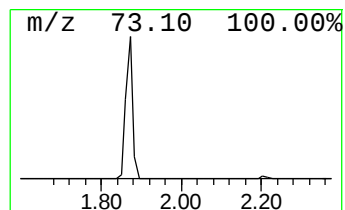
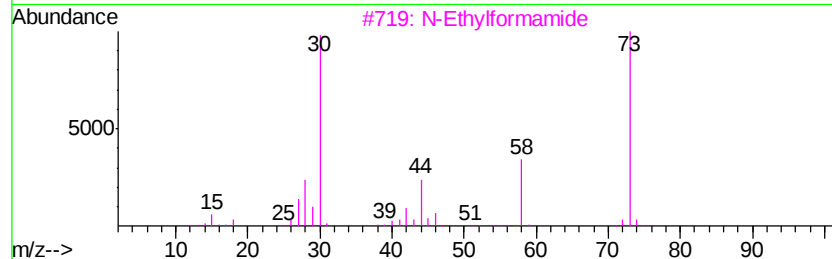
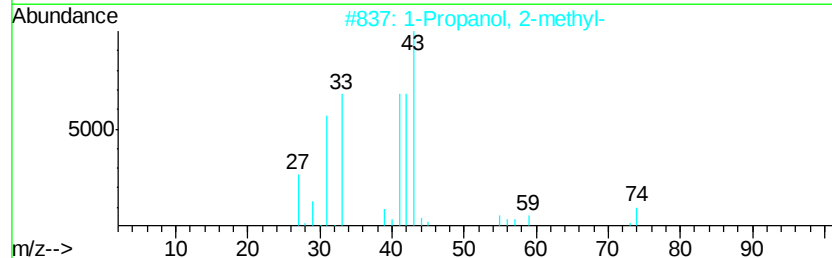
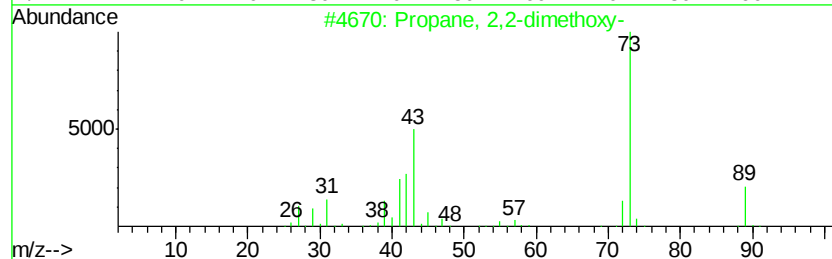
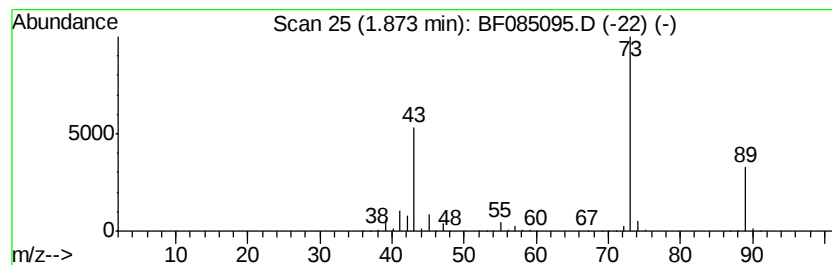
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF021716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Propane, 2,2-dimethoxy- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.87	104.75 ng	6195480	1,4-Dichlorobenzene-d4	6.99

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propane, 2,2-dimethoxy-	104	C5H12O2	000077-76-9	56
2		1-Propanol, 2-methyl-	74	C4H10O	000078-83-1	9
3		N-Ethylformamide	73	C3H7NO	000627-45-2	9
4		1,3-Diaminoquanidine	89	CH7N5	004364-78-7	9
5		2-Hydroxy-2-methylbutyric acid	118	C5H10O3	003739-30-8	9



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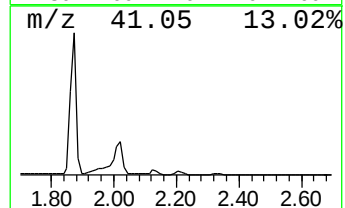
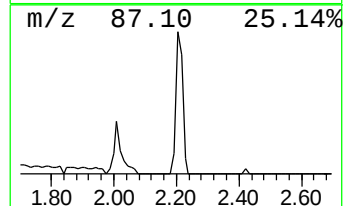
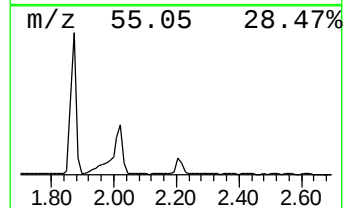
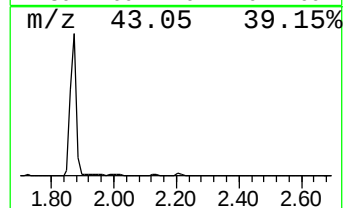
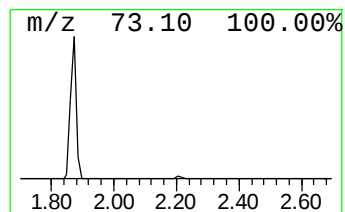
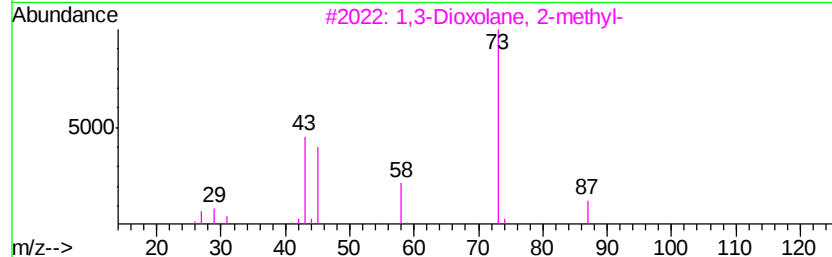
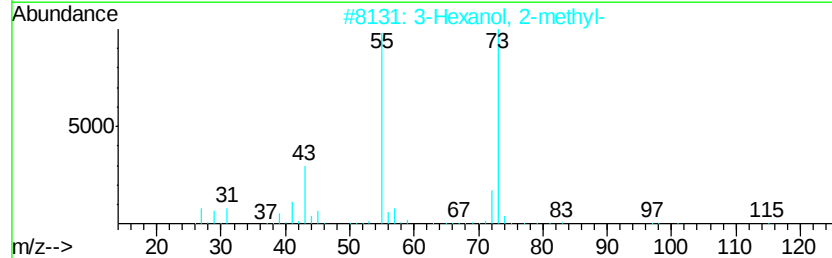
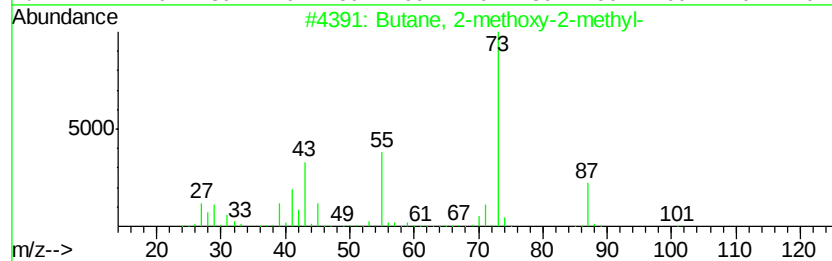
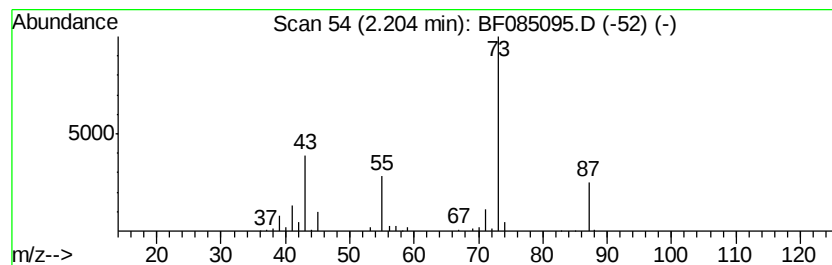
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Butane, 2-methoxy-2-methyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.20	2.76 ng	163182	1,4-Dichlorobenzene-d4	6.99

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	72
2		3-Hexanol, 2-methyl-	116	C7H16O	000617-29-8	25
3		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	25
4		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17
5		Silane, tetramethyl-	88	C4H12Si	000075-76-3	10



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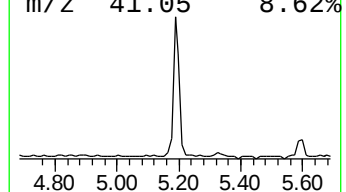
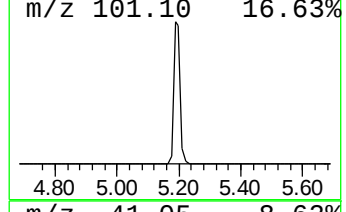
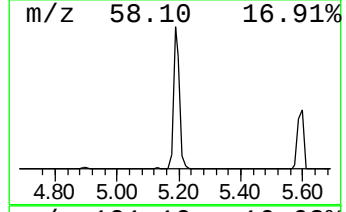
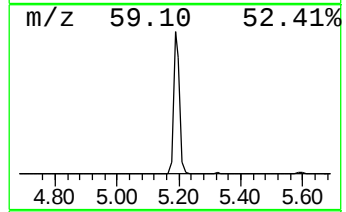
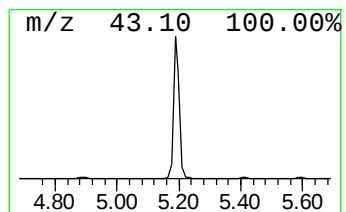
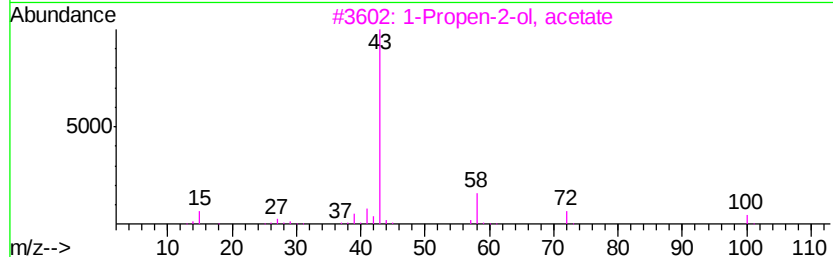
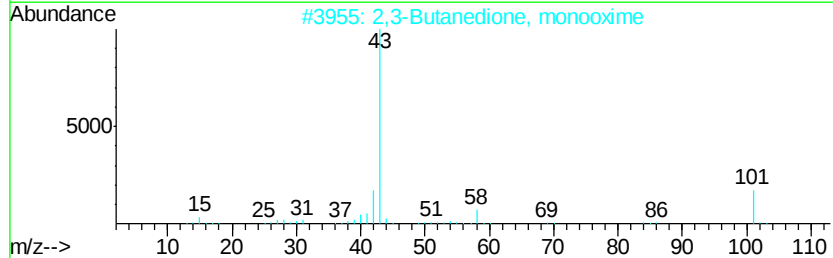
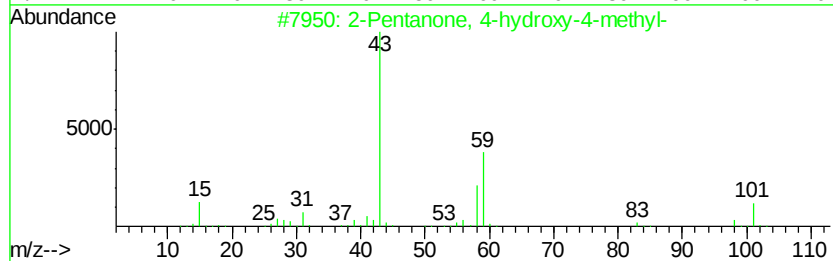
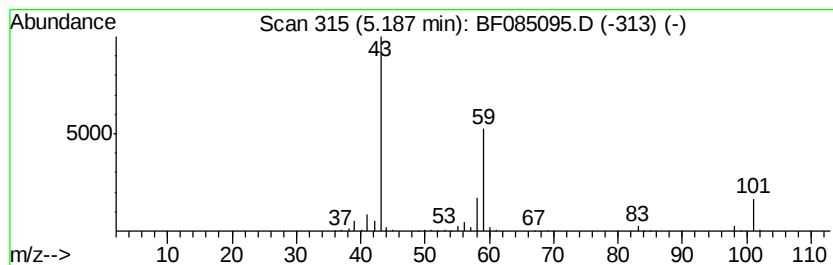
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.19	11.04 ng	652841	1,4-Dichlorobenzene-d4	6.99

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16
3		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
4		Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	9
5		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	9



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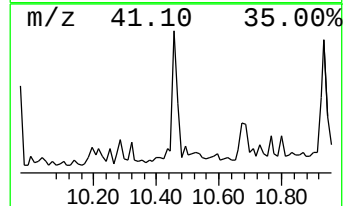
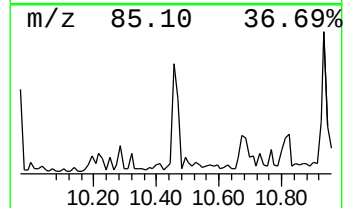
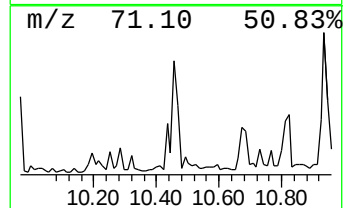
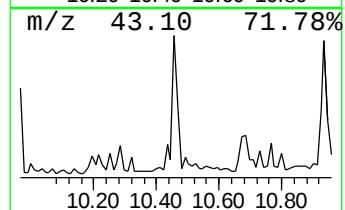
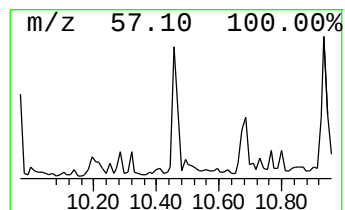
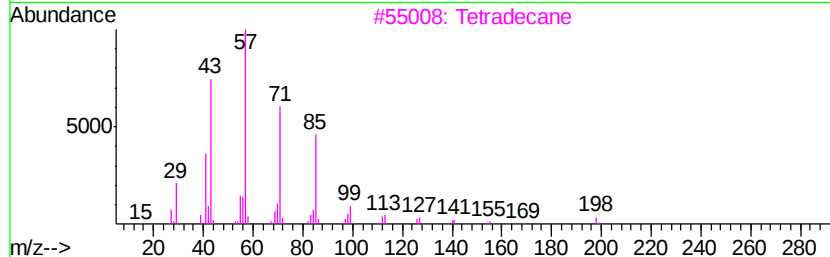
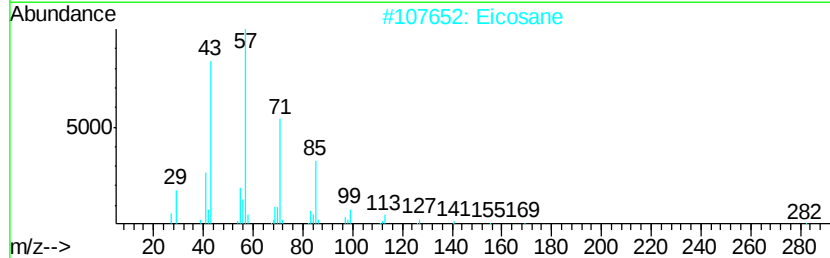
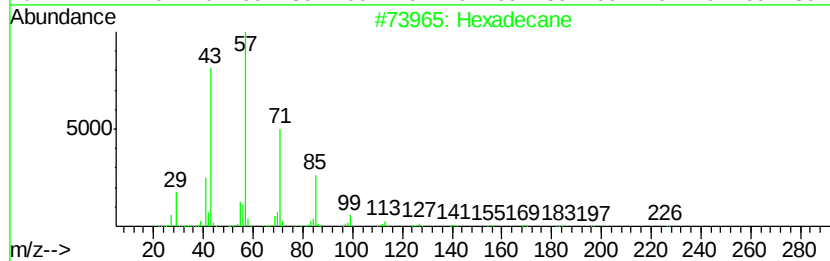
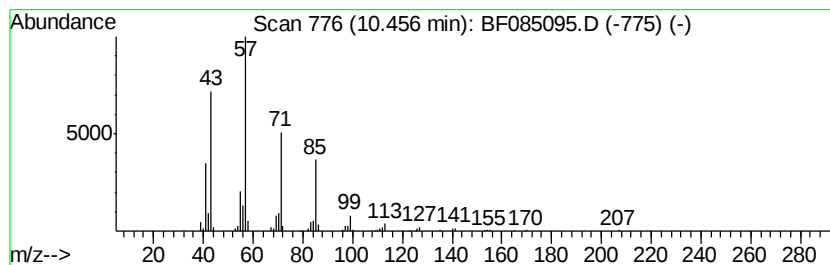
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Hexadecane Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.46	2.48 ng	220257	Acenaphthene-d10	10.03
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Hexadecane	226	C16H34	000544-76-3 90
2	Eicosane	282	C20H42	000112-95-8 87
3	Tetradecane	198	C14H30	000629-59-4 80
4	Nonane, 5-butyl-	184	C13H28	017312-63-9 80
5	Tridecane, 7-propyl-	226	C16H34	055045-09-5 80



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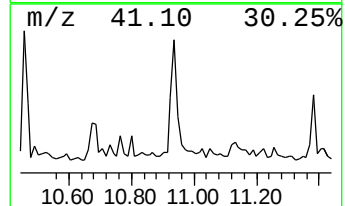
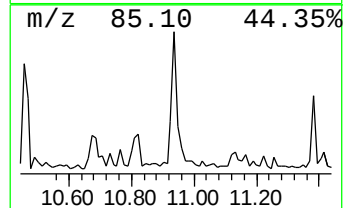
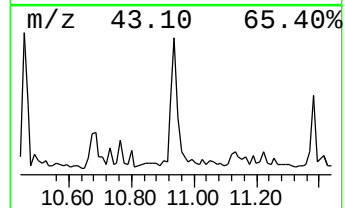
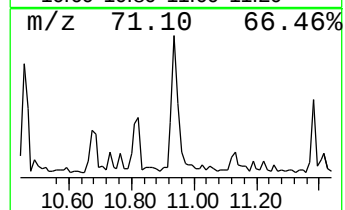
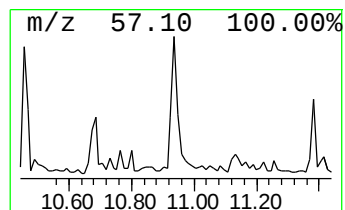
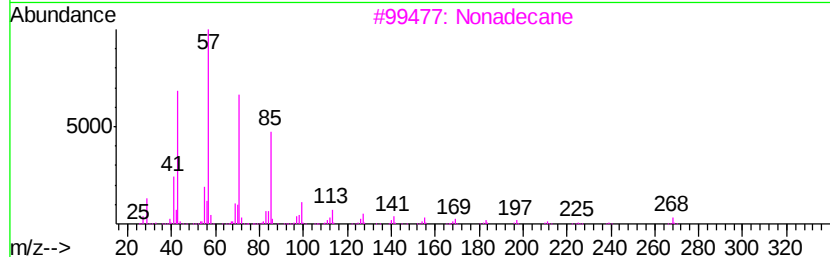
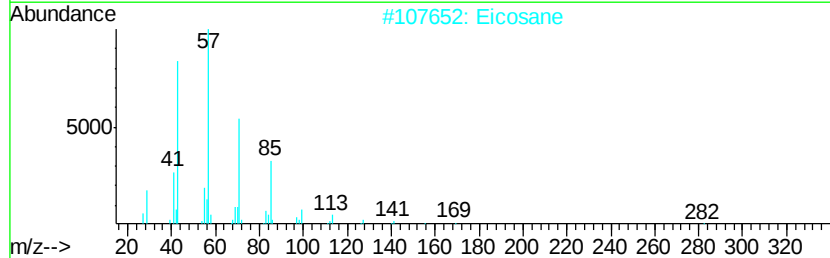
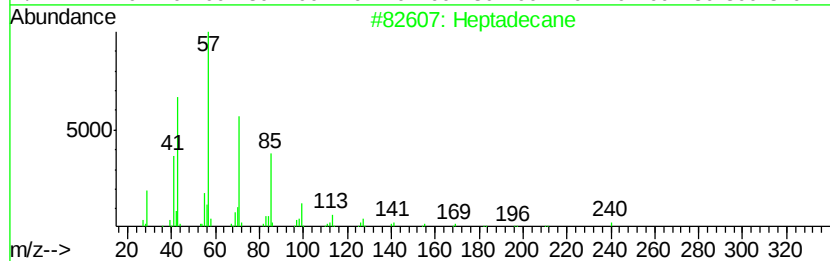
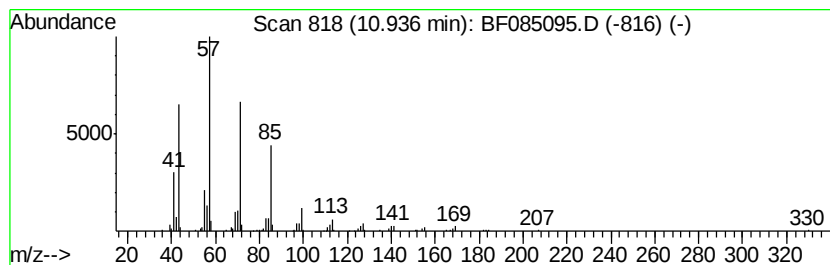
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Heptadecane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD		R.T.	
10.94	3.46 ng	320586	Phenanthrene-d10		11.52	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane		240	C17H36	000629-78-7	91
2	Eicosane		282	C20H42	000112-95-8	91
3	Nonadecane		268	C19H40	000629-92-5	91
4	Hexadecane		226	C16H34	000544-76-3	90
5	Tetradecane		198	C14H30	000629-59-4	87



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF021716\
Data File : BF085095.D
Acq On : 17 Feb 2016 19:20
Operator : SJ/IZ
Sample : H1443-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

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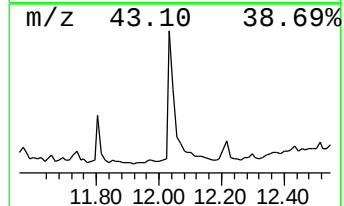
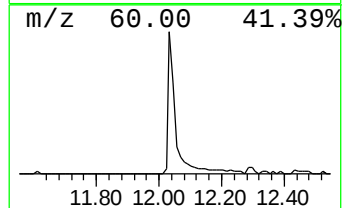
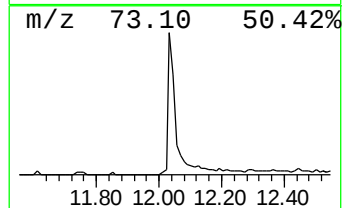
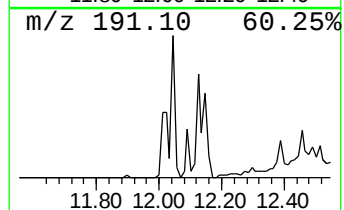
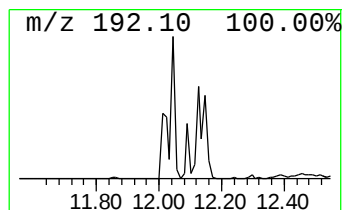
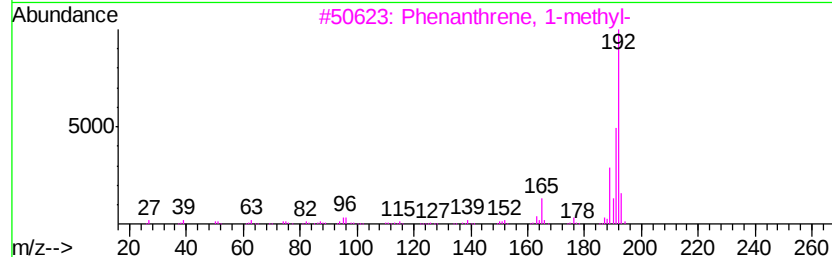
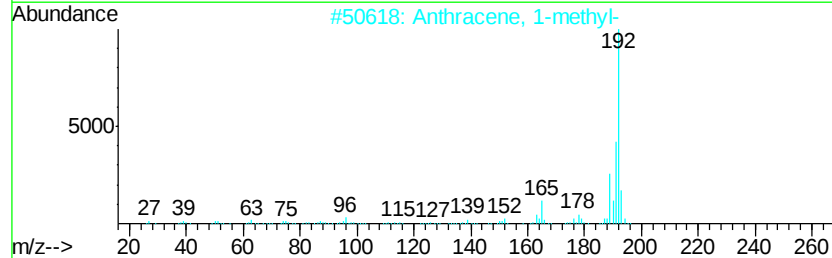
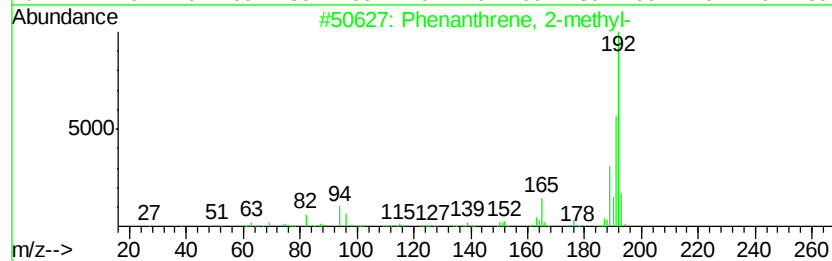
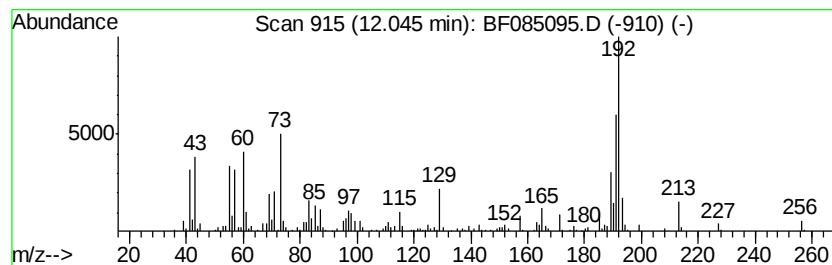
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Phenanthrene, 2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.05	5.05 ng	468196	Phenanthrene-d10	11.52

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
2		Anthracene, 1-methyl-	192	C15H12	000610-48-0	95
3		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95
4		Anthracene, 2-methyl-	192	C15H12	000613-12-7	93
5		Anthracene, 9-methyl-	192	C15H12	000779-02-2	89



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF021716\
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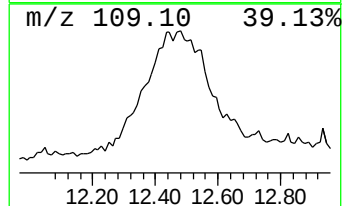
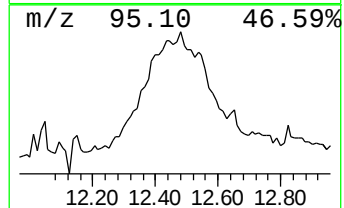
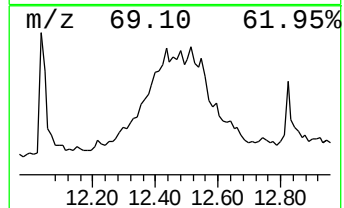
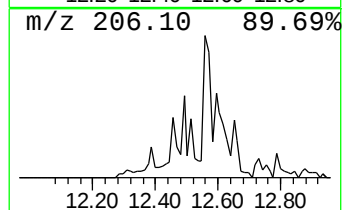
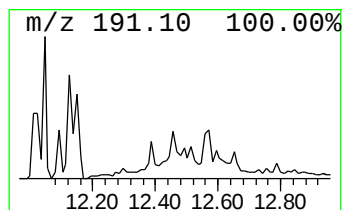
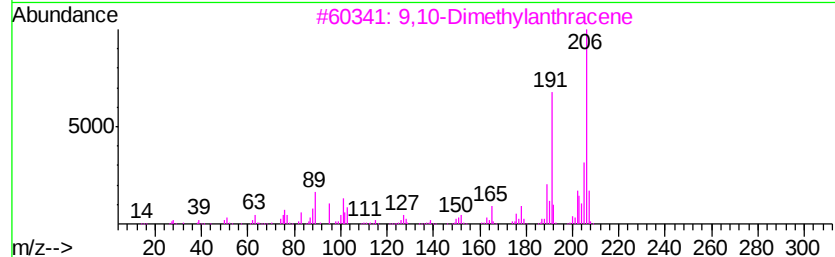
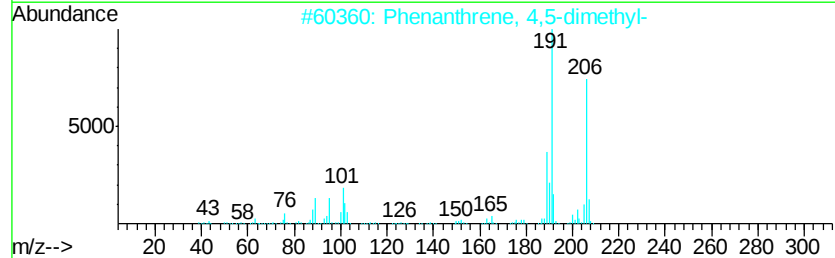
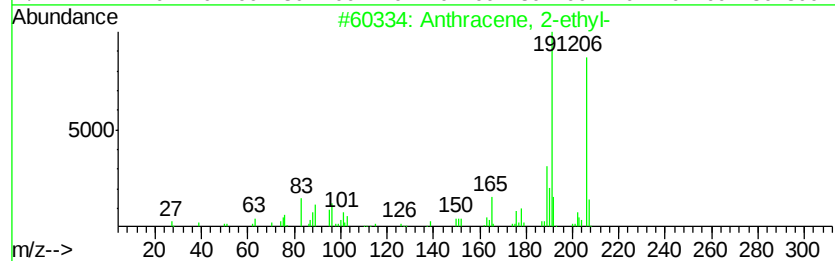
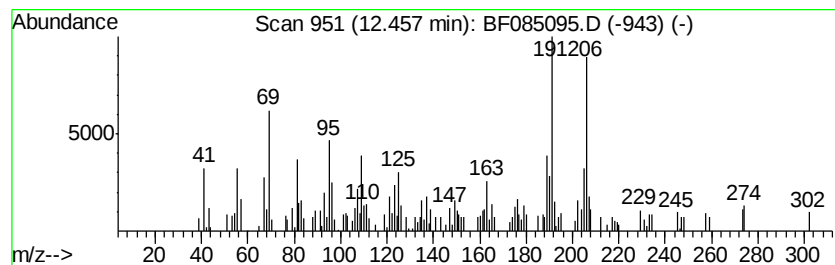
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Anthracene, 2-ethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.46	3.09 ng	286278	Phenanthrene-d10	11.52

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Anthracene, 2-ethyl-	206	C16H14	052251-71-5	64
2		Phenanthrene, 4,5-dimethyl-	206	C16H14	003674-69-9	64
3		9,10-Dimethylanthracene	206	C16H14	000781-43-1	60
4		1H-Inden-1-one, 2,3-dihydro-5,6-...	206	C12H14O3	004082-25-1	45
5		Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	41



Data Path : Z:\HPCHEM1\BNA F\DATA\BF021716\
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ClientSampleId :
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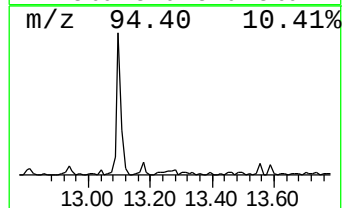
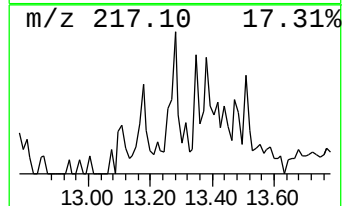
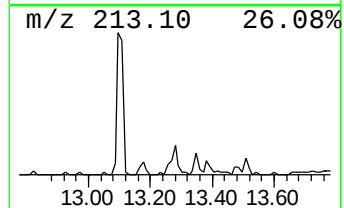
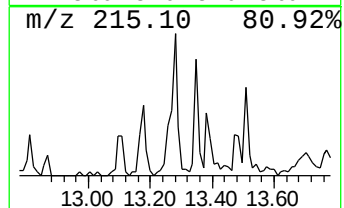
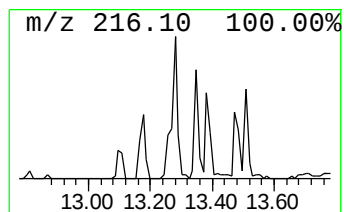
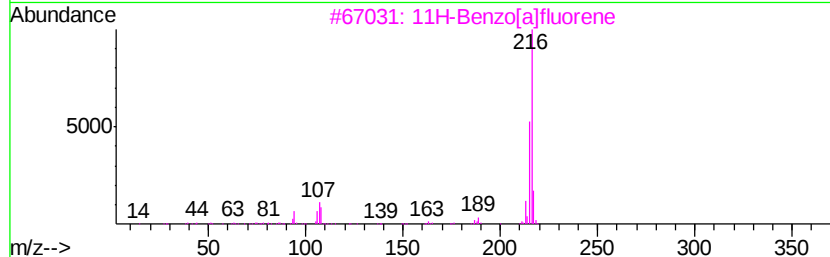
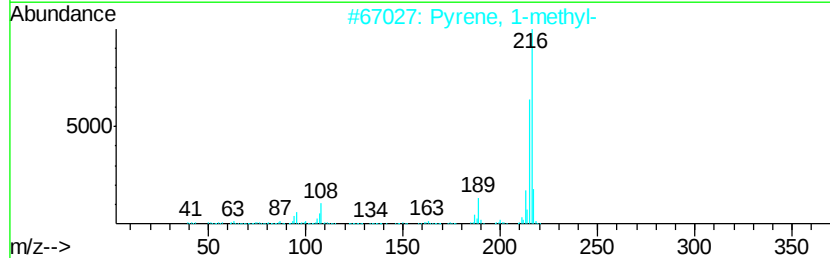
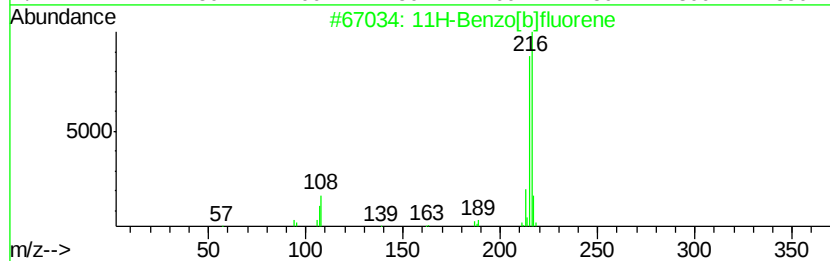
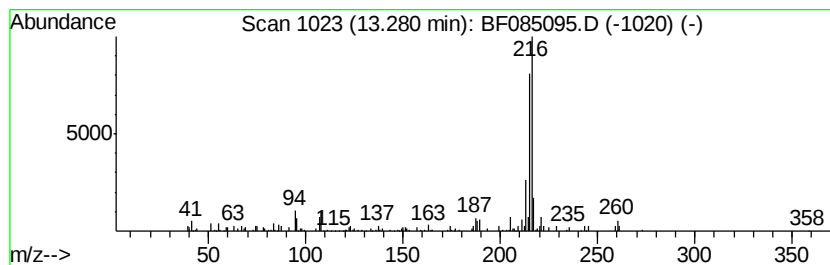
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 11H-Benzo[b]fluorene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.28	3.11 ng	254714	Chrysene-d12	14.15

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	90
2		Pyrene, 1-methyl-	216	C17H12	002381-21-7	87
3		11H-Benzo[a]fluorene	216	C17H12	000238-84-6	87
4		Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	81
5		.beta.-Carboline, 1,2,3,4-tetra...	260	C14H16N2O3	1000124-14-2	78



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF021716\
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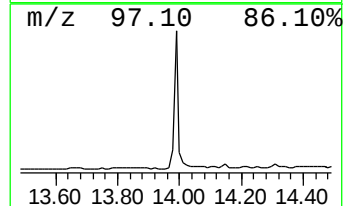
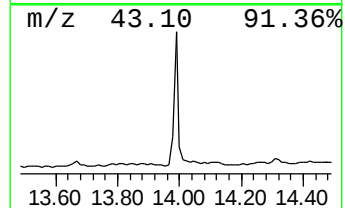
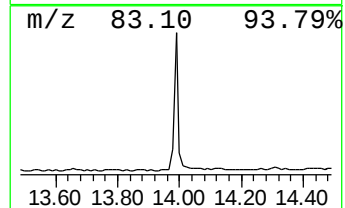
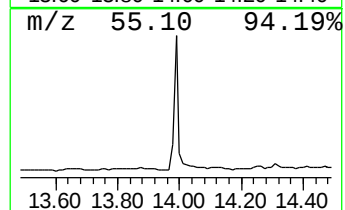
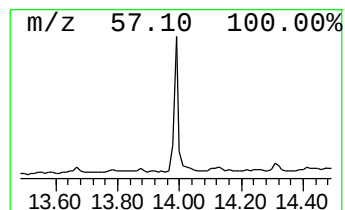
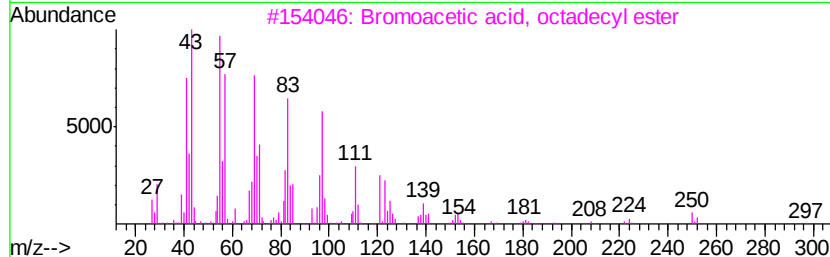
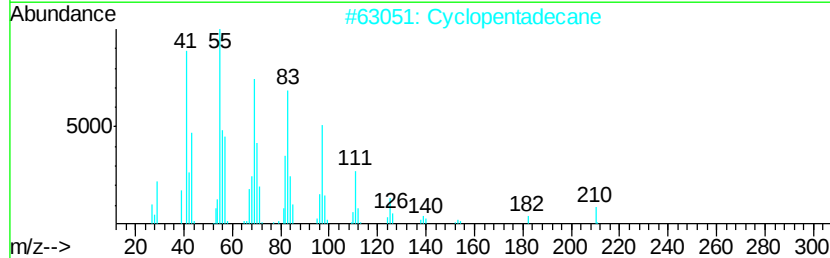
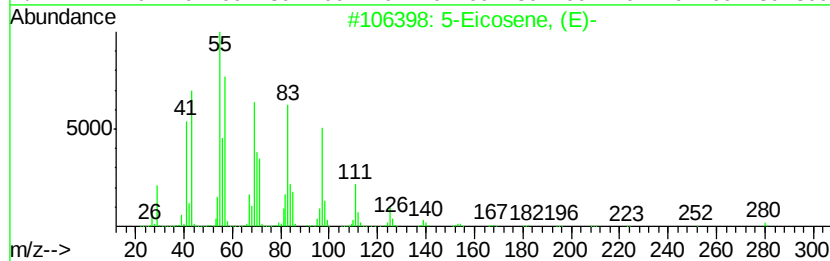
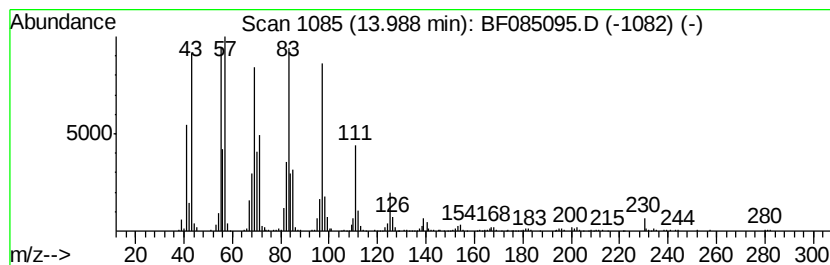
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 5-Eicosene, (E)- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.99	9.64 ng	790796	Chrysene-d12	14.15

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			5-Eicosene, (E)-	280	C20H40	074685-30-6	99
2			Cyclopentadecane	210	C15H30	000295-48-7	93
3			Bromoacetic acid, octadecyl ester	390	C20H39BrO2	018992-03-5	91
4			Trifluoroacetic acid, n-octadecyl...	366	C20H37F3O2	079392-43-1	91
5			1-Nonadecene	266	C19H38	018435-45-5	91



Data Path : Z:\HPCHEM1\BNA F\DATA\BF021716\
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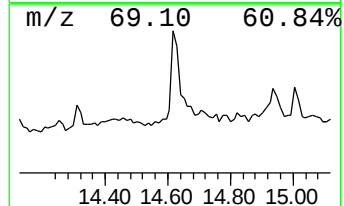
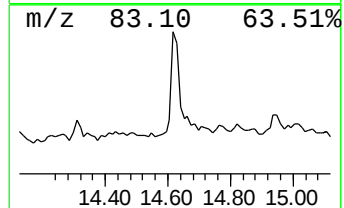
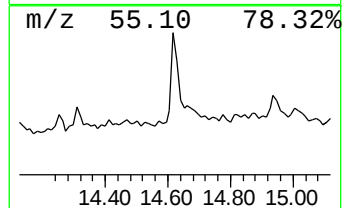
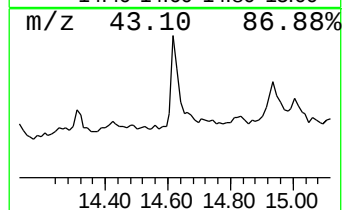
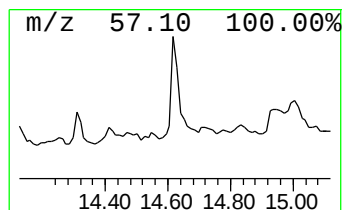
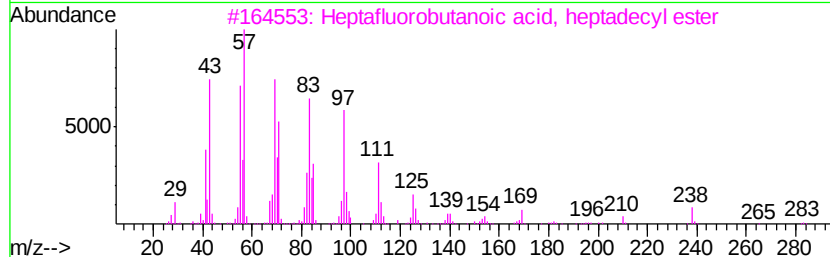
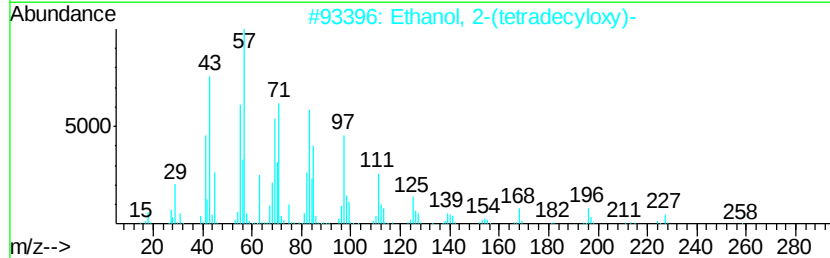
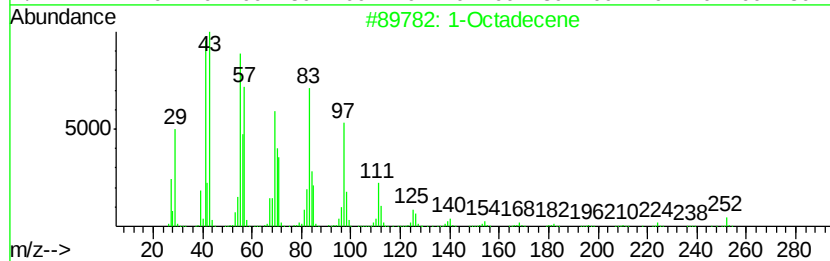
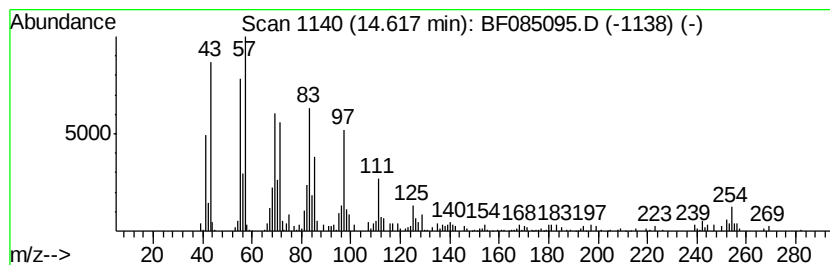
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 1-Octadecene Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.62	3.01 ng	246649	Chrysene-d12	14.15
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1-Octadecene	252 C18H36	000112-88-9	95
2	Ethanol, 2-(tetradecyloxy)-	258 C16H34O2	002136-70-1	87
3	Heptafluorobutanoic acid, heptad...	452 C21H35F7O2	1000282-97-3	87
4	Trifluoroacetic acid, n-octadecy...	366 C20H37F3O2	079392-43-1	83
5	Heptafluorobutyric acid, n-penta...	424 C19H31F7O2	1000216-79-3	76



Data Path : Z:\HPCHEM1\BNA F\DATA\BF021716\
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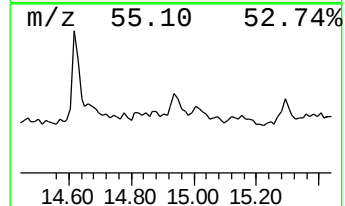
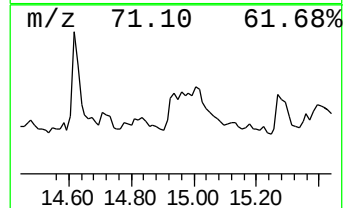
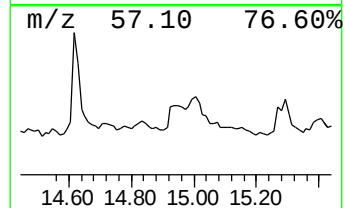
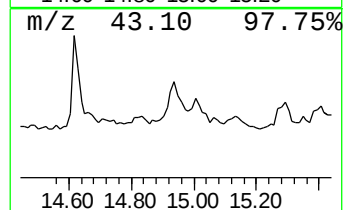
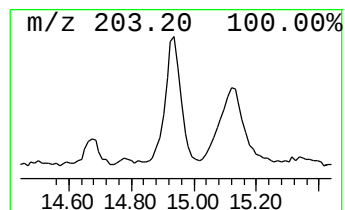
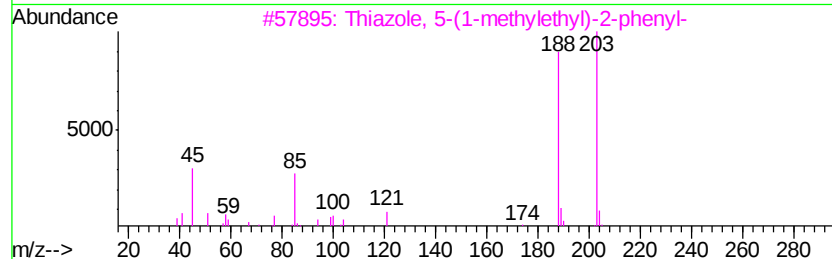
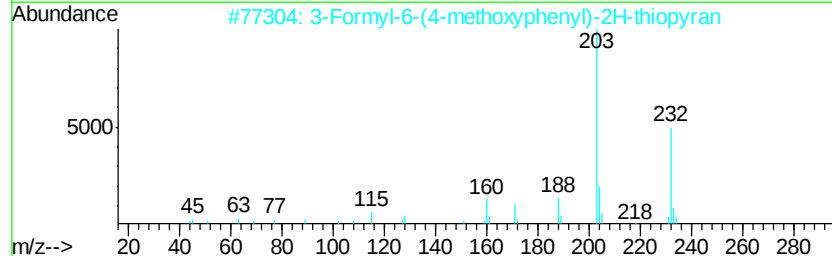
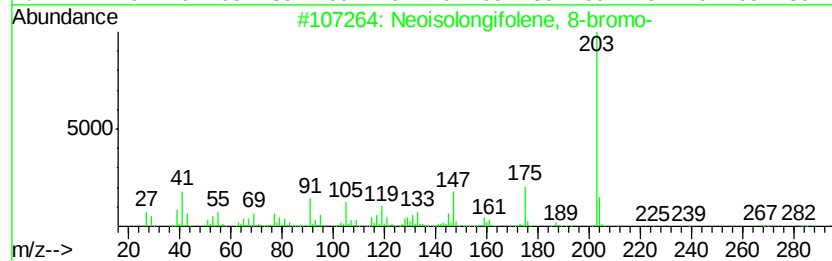
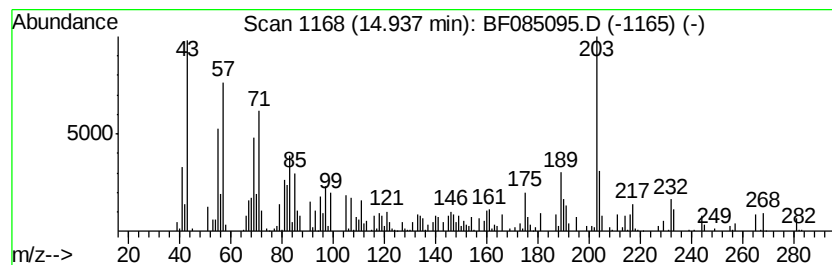
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 unknown14.94 Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.94	2.72 ng	152896	Perylene-d12	15.62

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Neoisonlongifolene, 8-bromo-	282	C15H23Br	1000151-64-8	41
2		3-Formyl-6-(4-methoxyphenyl)-2H-...	232	C13H12O2S	037750-09-7	38
3		Thiazole, 5-(1-methylethyl)-2-ph...	203	C12H13NS	039199-35-4	37
4		Phenol, 2-(4-ethoxycarbonylthiaz...	249	C12H11NO3S	027383-13-7	35
5		Pyrano[4,3-b]pyrane-3-carbonitri...	280	C16H12N2O3	194282-72-9	27



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF021716\
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Misc :
ALS Vial : 15 Sample Multiplier: 1

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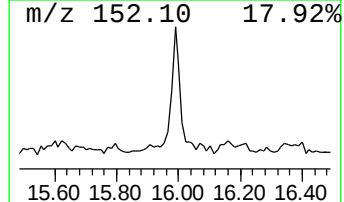
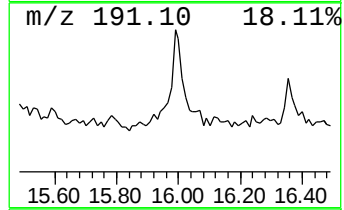
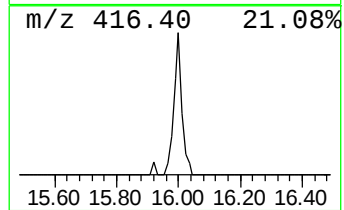
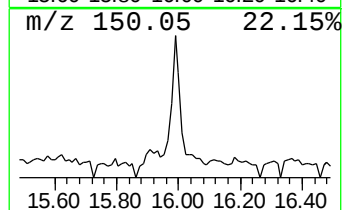
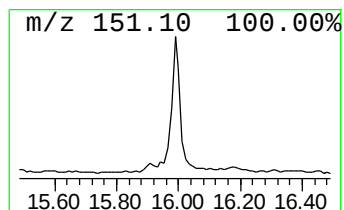
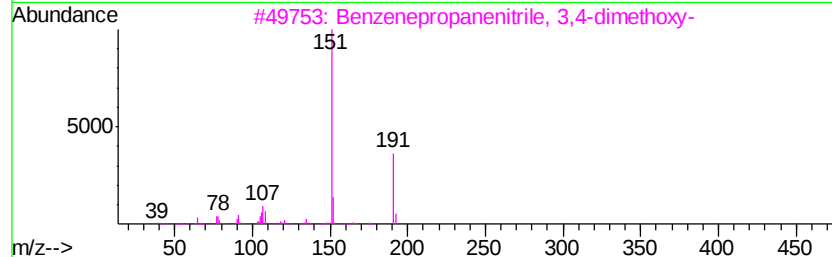
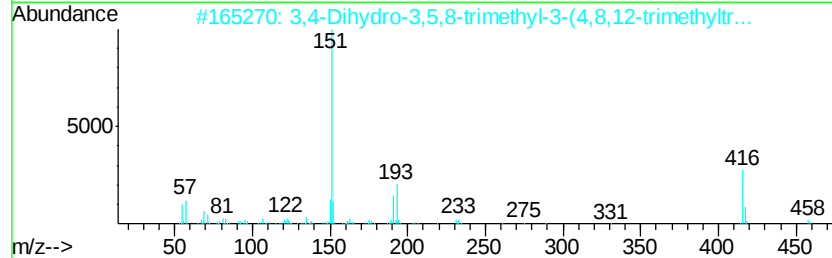
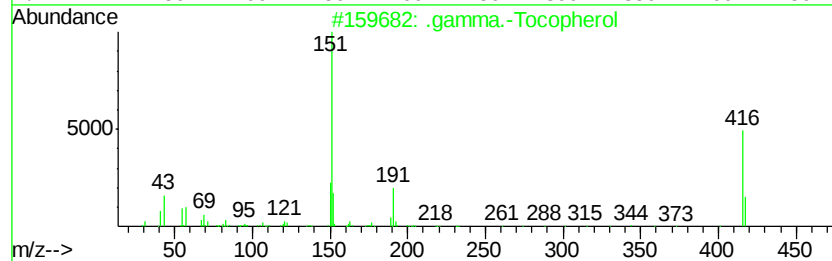
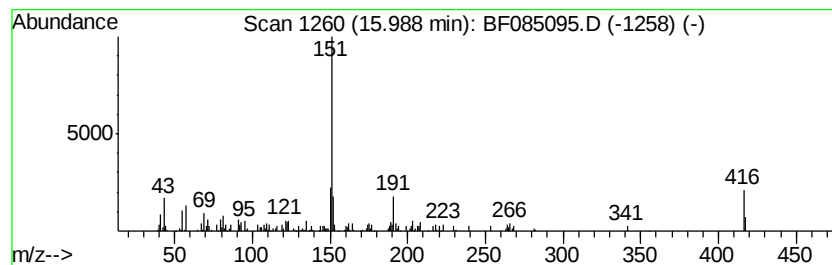
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 .gamma.-Tocopherol Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.99	3.80 ng	214003	Perylene-d12	15.62

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	.gamma.-Tocopherol	416	C28H48O2	007616-22-0	93
2		3,4-Dihydro-3,5,8-trimethyl-3-(4...	458	C30H50O3	1000105-71-9	72
3		Benzenepropanenitrile, 3,4-dimet...	191	C11H13N02	049621-56-9	58
4		3-Amino-4-methylbenzoic acid	151	C8H9NO2	002458-12-0	40
5		Dimethoxydimethylgermanium	166	C4H12GeO2	004531-27-5	38



Data Path : Z:\HPCHEM1\BNA F\DATA\BF021716\
Data File : BF085095.D
Acq On : 17 Feb 2016 19:20
Operator : SJ/IZ
Sample : H1443-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
EP004

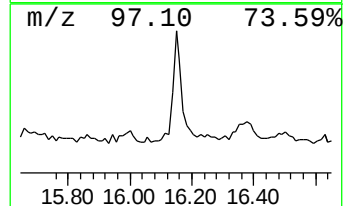
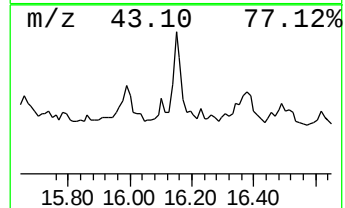
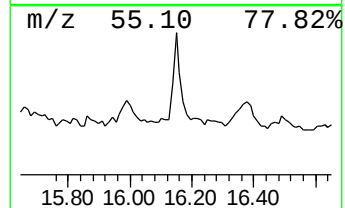
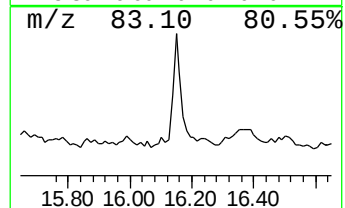
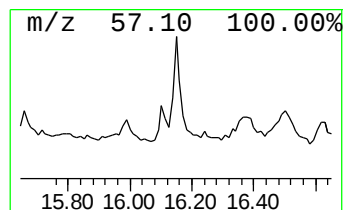
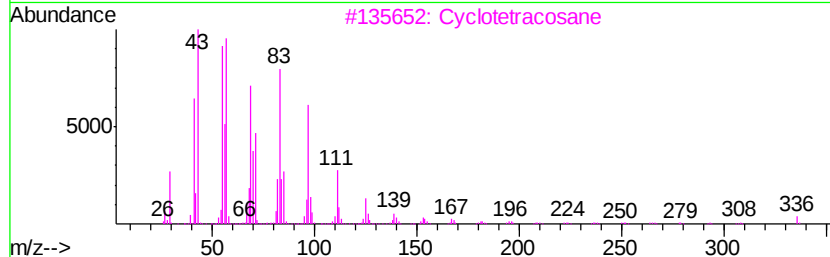
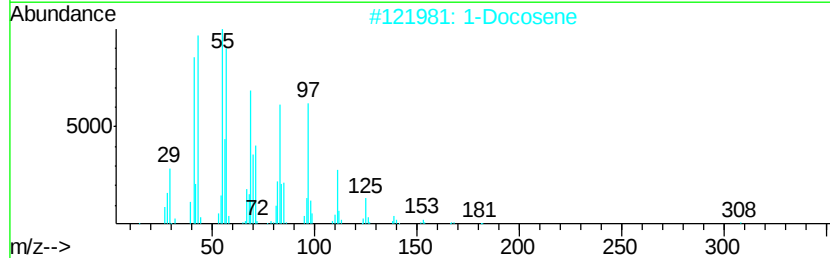
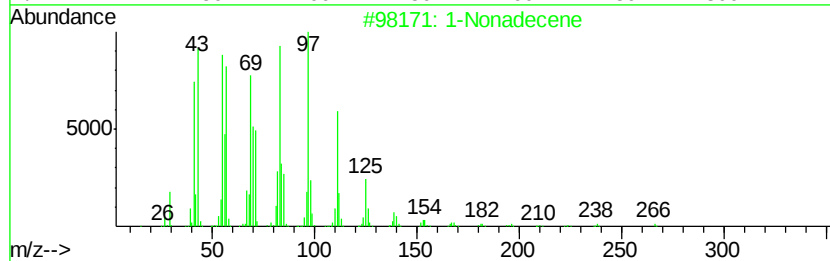
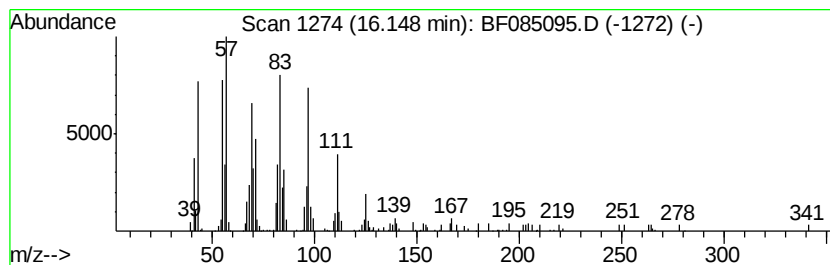
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 1-Nonadecene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.15	3.90 ng	219493	Perylene-d12	15.62

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Nonadecene	266	C19H38	018435-45-5	95
2		1-Docosene	308	C22H44	001599-67-3	86
3		Cyclotetracosane	336	C24H48	000297-03-0	83
4		Heptafluorobutanoic acid, heptad...	452	C21H35F7O2	1000282-97-3	74
5		Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	70



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF021716\
Data File : BF085095.D
Acq On : 17 Feb 2016 19:20
Operator : SJ/IZ
Sample : H1443-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
EP004

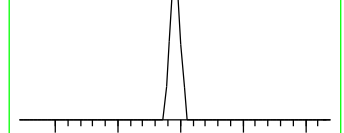
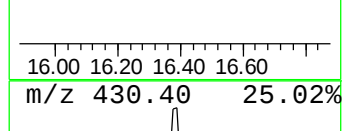
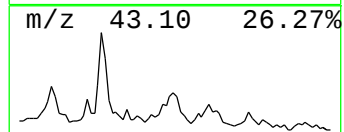
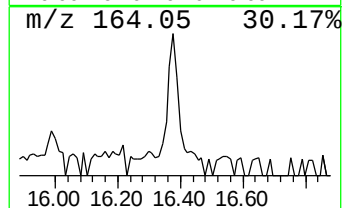
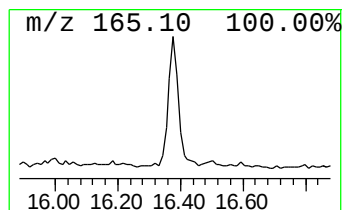
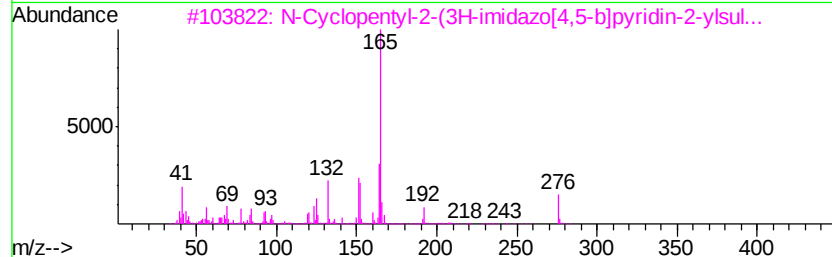
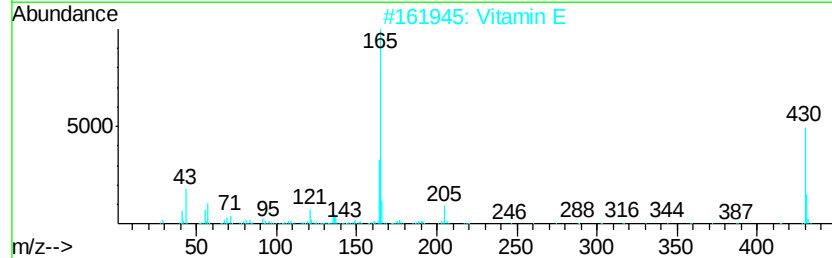
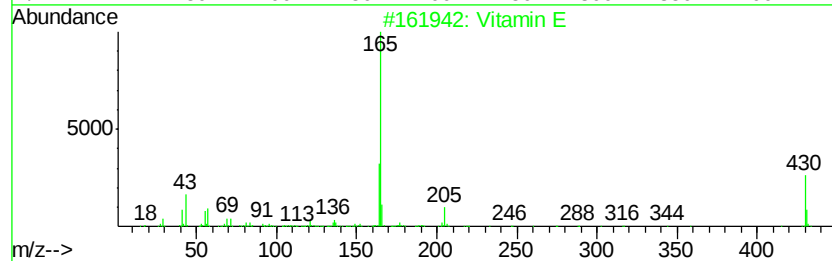
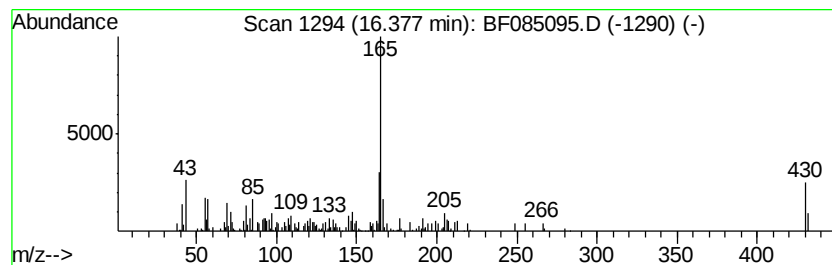
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 Vitamin E Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.38	3.51 ng	197777	Perylene-d12	15.62

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Vitamin E	430	C29H50O2	010191-41-0	87
2		Vitamin E	430	C29H50O2	000059-02-9	87
3		N-Cyclopentyl-2-(3H-imidazo[4,5-b]pyridin-2-ylsul...	276	C13H16N4OS	1000295-07-4	50
4		4-Amino-3-methoxypyrazolo[3,4-d]pyrimidin-2(1H)-one	165	C6H7N5O	111375-22-5	47
5		6H-Purin-6-one, 1,7-dihydro-2-(methylthio)-	165	C6H7N5O	010030-78-1	43



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF021716\
 Data File : BF085095.D
 Acq On : 17 Feb 2016 19:20
 Operator : SJ/IZ
 Sample : H1443-01
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EP004

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF021716.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
Propane, 2,2-dime...	1.87	104.8	ng	6195480	1	6.99	1182890	20.0
Butane, 2-methoxy...	2.20	2.8	ng	163182	1	6.99	1182890	20.0
2-Pentanone, 4-hy...	5.19	11.0	ng	652841	1	6.99	1182890	20.0
Hexadecane	10.46	2.5	ng	220257	3	10.03	1773470	20.0
Heptadecane	10.94	3.5	ng	320586	4	11.52	1855650	20.0
Phenanthrene, 2-m...	12.05	5.0	ng	468196	4	11.52	1855650	20.0
Anthracene, 2-ethyl-	12.46	3.1	ng	286278	4	11.52	1855650	20.0
11H-Benzo[b]fluorene	13.28	3.1	ng	254714	5	14.15	1640350	20.0
5-Eicosene, (E)-	13.99	9.6	ng	790796	5	14.15	1640350	20.0
1-Octadecene	14.62	3.0	ng	246649	5	14.15	1640350	20.0
unknown14.94	14.94	2.7	ng	152896	6	15.62	1125430	20.0
.gamma.-Tocopherol	15.99	3.8	ng	214003	6	15.62	1125430	20.0
1-Nonadecene	16.15	3.9	ng	219493	6	15.62	1125430	20.0
Vitamin E	16.38	3.5	ng	197777	6	15.62	1125430	20.0