

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042915\
 Data File : BF078837.D
 Acq On : 30 Apr 2015 4:41
 Operator : TP/IZ
 Sample : G2021-11
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PS-25-0-2

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-BF042815.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	1.290	6	9	11	rVB	1813159	2589756	52.11%	6.720%
2	1.530	27	30	33	rVB	4586681	4970047	100.00%	12.897%
3	1.678	40	43	46	rVB	1619892	2162327	43.51%	5.611%
4	1.781	51	52	56	rBV	182291	284920	5.73%	0.739%
5	1.861	56	59	62	rVB	409570	580727	11.68%	1.507%
6	1.918	62	64	89	rVB	2681135	4604256	92.64%	11.948%
7	5.153	345	347	349	rBV	57558	74173	1.49%	0.192%
8	5.187	349	350	356	rVB	208436	220899	4.44%	0.573%
9	5.679	390	393	395	rBV	1286235	1370302	27.57%	3.556%
10	6.925	500	502	505	rBV	1776009	1584511	31.88%	4.112%
11	7.085	514	516	519	rBV	1126724	1530073	30.79%	3.970%
12	7.382	540	542	544	rBV	535528	476589	9.59%	1.237%
13	7.565	556	558	561	rVB	949161	1175095	23.64%	3.049%
14	8.067	600	602	605	rBV	888252	890239	17.91%	2.310%
15	8.959	678	680	683	rBV	797194	698046	14.05%	1.811%
16	10.308	795	798	800	rBV	2306440	2032796	40.90%	5.275%
17	10.811	840	842	845	rBV	63200	75971	1.53%	0.197%
18	11.096	864	867	868	rVB	36466	52569	1.06%	0.136%
19	11.142	868	871	873	rBV	705256	766950	15.43%	1.990%
20	11.714	919	921	923	rVB	99060	89157	1.79%	0.231%
21	11.988	941	945	947	rBV	53930	79567	1.60%	0.206%
22	12.114	954	956	959	rVB	1018076	1101414	22.16%	2.858%
23	12.308	971	973	978	rBV	188223	297925	5.99%	0.773%
24	12.868	1020	1022	1024	rBV	236308	235660	4.74%	0.612%
25	12.902	1024	1025	1030	rVB	95613	103428	2.08%	0.268%
26	12.982	1030	1032	1036	rBV	766624	906466	18.24%	2.352%
27	13.085	1039	1041	1045	rVV2	34247	63560	1.28%	0.165%
28	13.348	1062	1064	1065	rBV	42800	52682	1.06%	0.137%
29	13.405	1066	1069	1071	rVV	244390	234736	4.72%	0.609%
30	13.611	1083	1087	1089	rBV3	48046	117196	2.36%	0.304%
31	13.691	1092	1094	1096	rBV	61174	72057	1.45%	0.187%
32	13.919	1112	1114	1115	rBV	166927	226575	4.56%	0.588%
33	13.988	1118	1120	1125	rBV5	42262	80313	1.62%	0.208%
34	14.114	1128	1131	1134	rVV4	31717	64267	1.29%	0.167%

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Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF042815.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	14.262	1143	1144	1146	rBV	52191	68431	1.38%	0.178%
36	14.400	1154	1156	1158	rBV	190653	197033	3.96%	0.511%
37	14.491	1162	1164	1166	rBV	157813	178896	3.60%	0.464%
38	14.537	1166	1168	1170	rBV	38058	62774	1.26%	0.163%
39	14.571	1170	1171	1174	rBV3	44208	80850	1.63%	0.210%
40	14.674	1178	1180	1181	rBV	58937	78759	1.58%	0.204%
41	14.777	1187	1189	1191	rVB	167389	184747	3.72%	0.479%
42	14.845	1192	1195	1196	rBV2	128661	204829	4.12%	0.532%
43	14.868	1196	1197	1199	rVB	169717	144879	2.92%	0.376%
44	14.982	1204	1207	1209	rBV	2319953	2379156	47.87%	6.174%
45	15.017	1209	1210	1216	rVB4	68512	177156	3.56%	0.460%
46	15.314	1234	1236	1238	rVB3	107971	119119	2.40%	0.309%
47	15.371	1238	1241	1244	rBV3	75011	173060	3.48%	0.449%
48	15.474	1248	1250	1252	rBV	106810	145480	2.93%	0.378%
49	15.748	1272	1274	1275	rBV	141804	137249	2.76%	0.356%
50	16.148	1305	1309	1315	rBV3	210673	729082	14.67%	1.892%
51	16.285	1318	1321	1323	rBV	709451	1289881	25.95%	3.347%
52	17.086	1388	1391	1394	rBV	264948	522031	10.50%	1.355%
53	18.011	1467	1472	1475	rBV	869860	1531250	30.81%	3.973%
54	19.223	1576	1578	1582	rVB	158464	267149	5.38%	0.693%

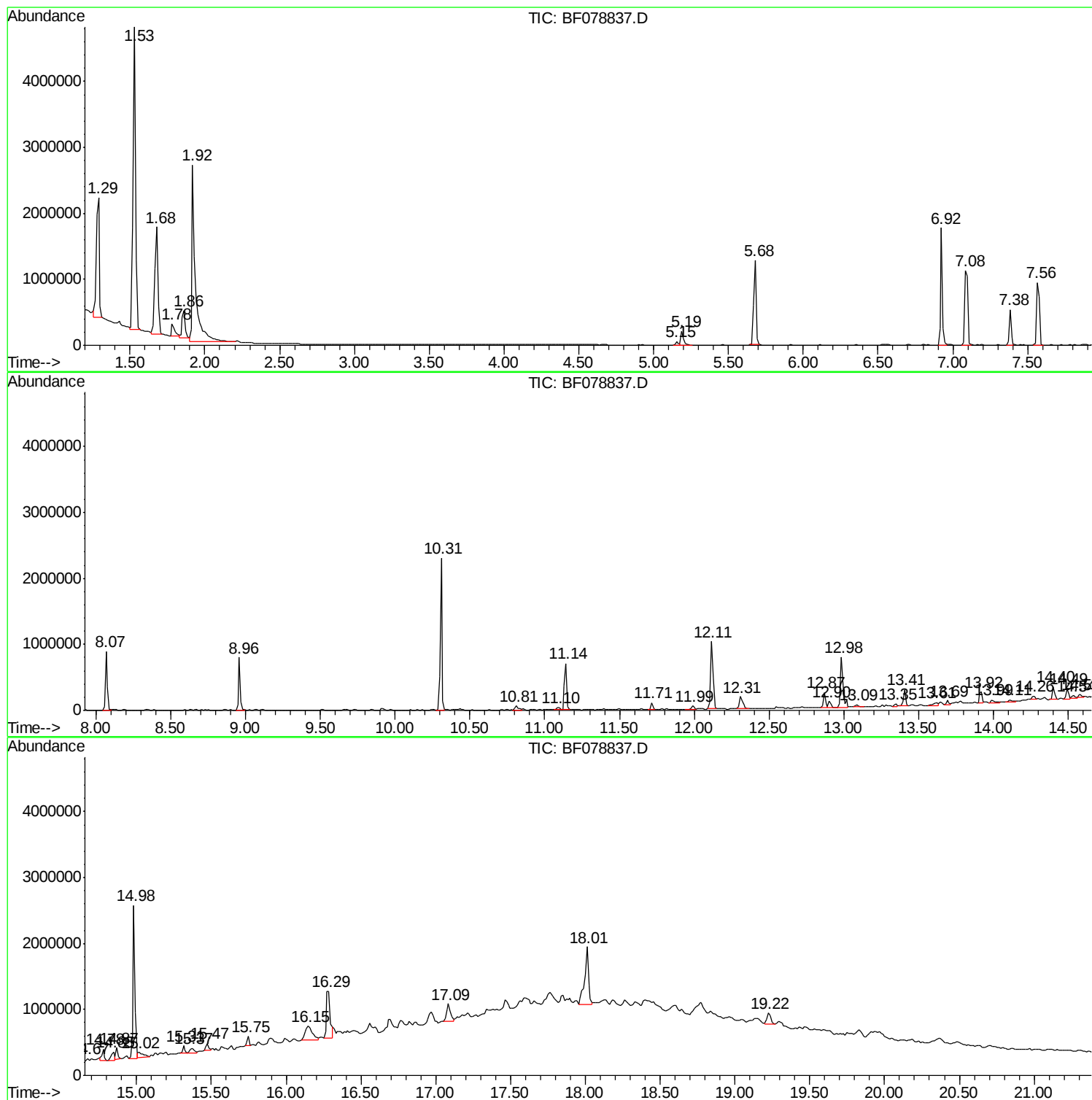
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF042815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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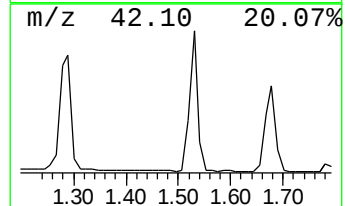
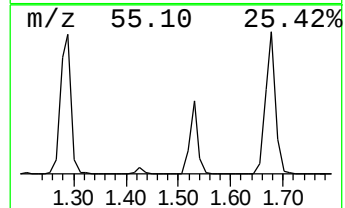
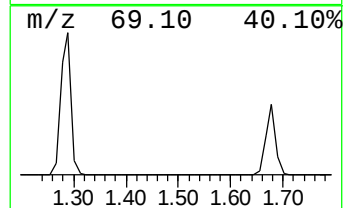
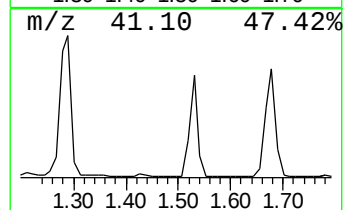
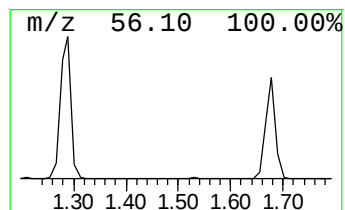
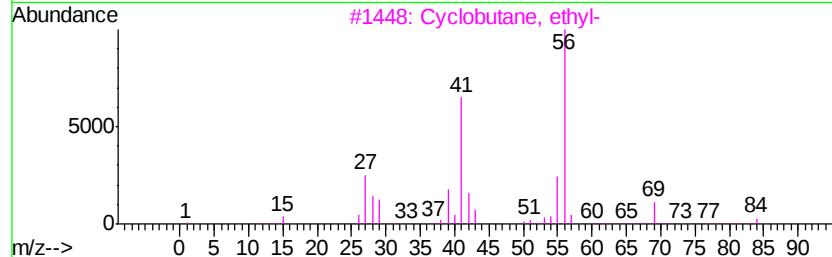
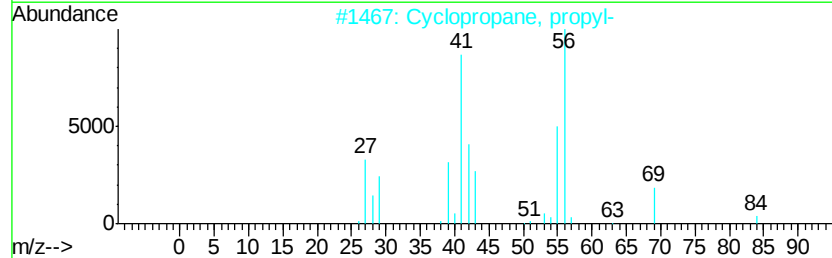
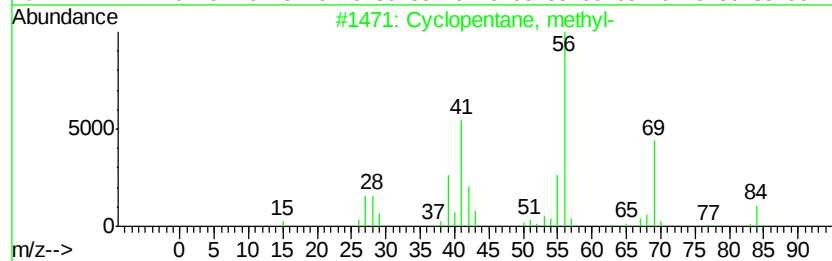
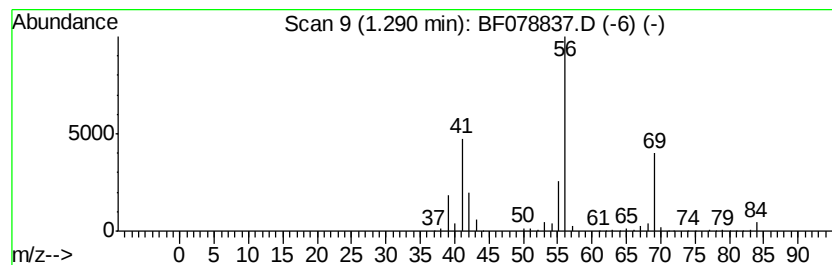
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF042815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Cyclopentane, methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.29	108.68 ng	2589760	1,4-Dichlorobenzene-d4	7.38

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentane, methyl-	84	C6H12	000096-37-7	90
2		Cyclopropane, propyl-	84	C6H12	002415-72-7	72
3		Cyclobutane, ethyl-	84	C6H12	004806-61-5	64
4		Cyclopentane, 1,1-dimethyl-	98	C7H14	001638-26-2	56
5		1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	39



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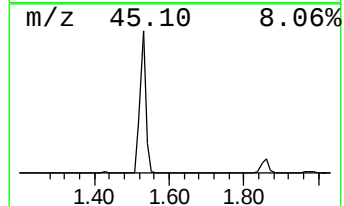
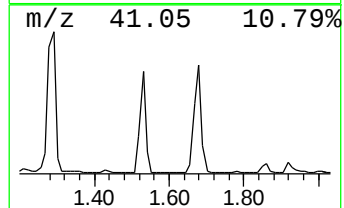
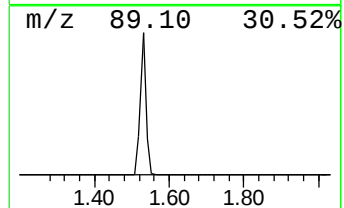
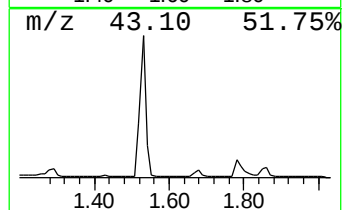
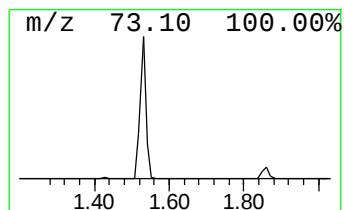
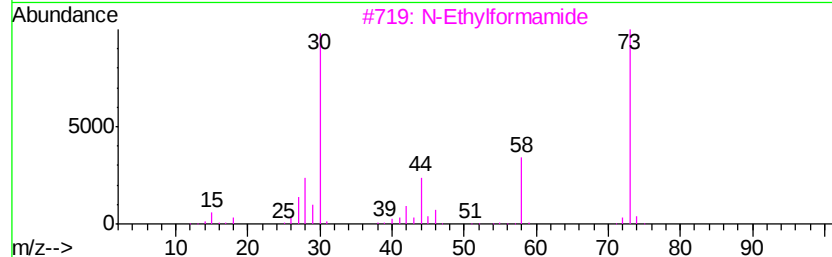
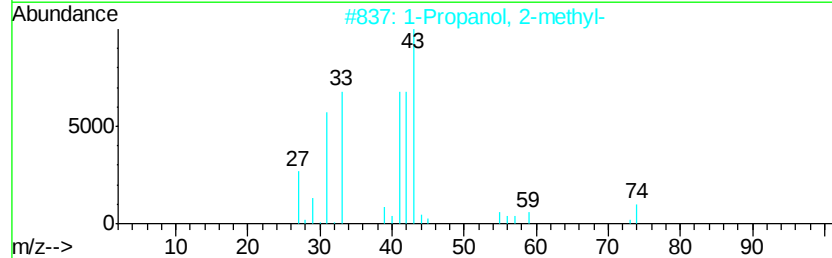
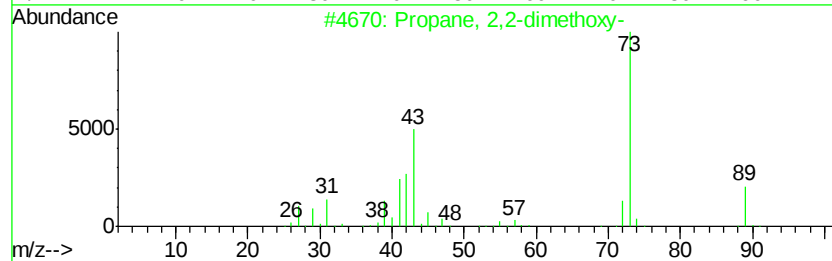
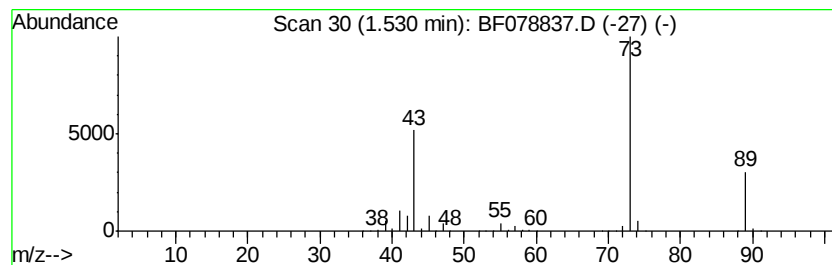
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.53	208.57 ng	4970050	1,4-Dichlorobenzene-d4	7.38

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propane, 2,2-dimethoxy-	104	C5H12O2	000077-76-9	72
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		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	9



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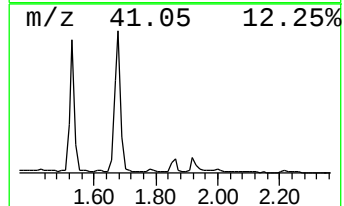
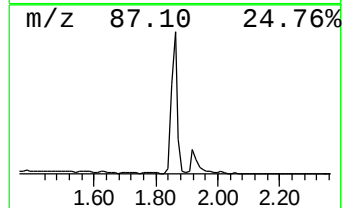
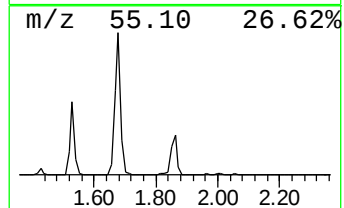
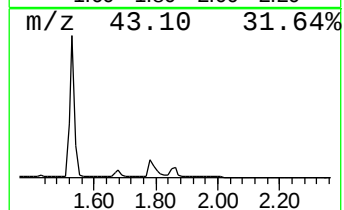
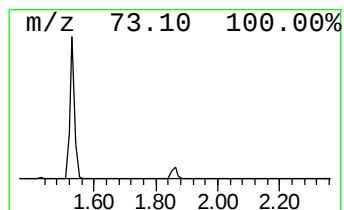
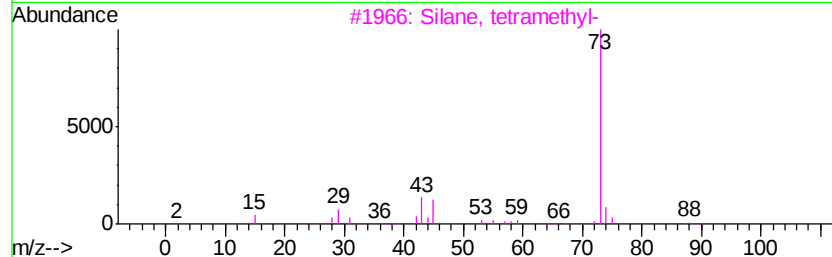
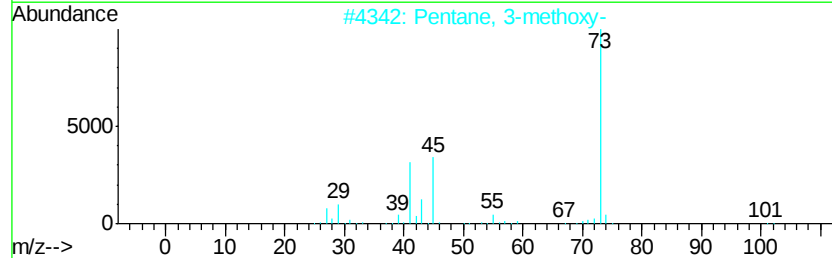
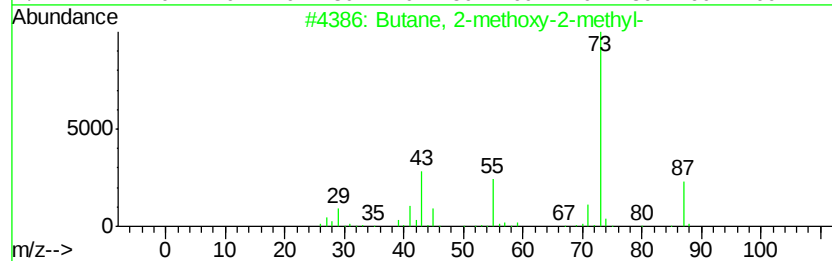
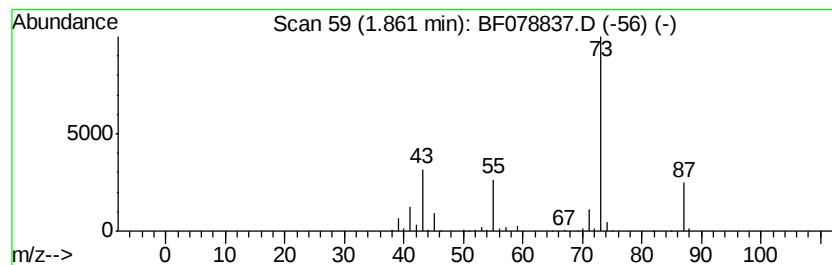
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Butane, 2-methoxy-2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.86	24.37 ng	580727	1,4-Dichlorobenzene-d4	7.38

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17
3		Silane, tetramethyl-	88	C4H12Si	000075-76-3	12
4		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9
5		N-Ethylformamide	73	C3H7NO	000627-45-2	9



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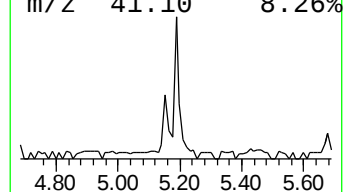
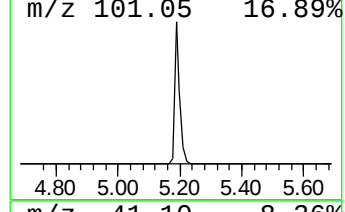
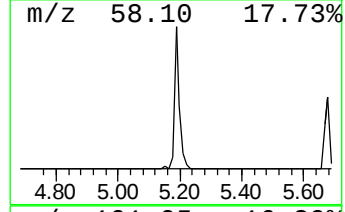
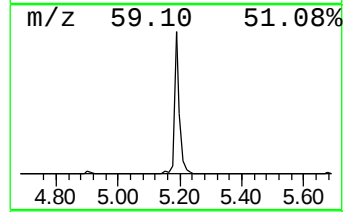
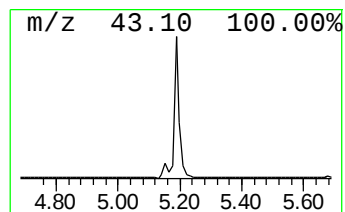
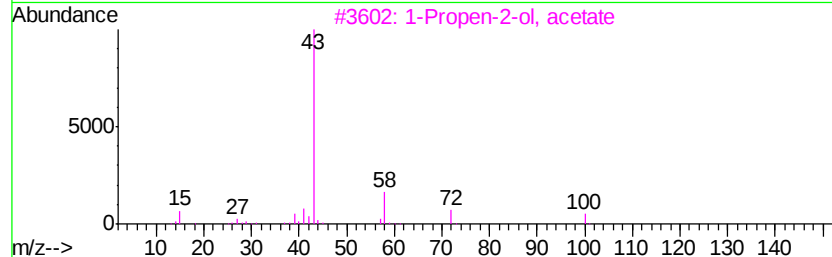
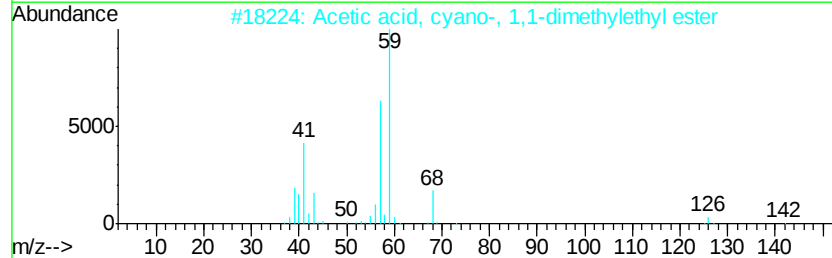
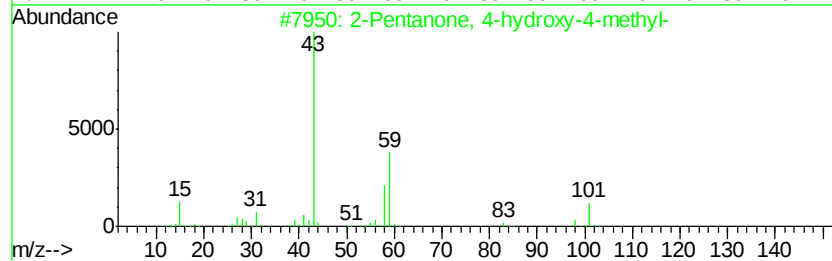
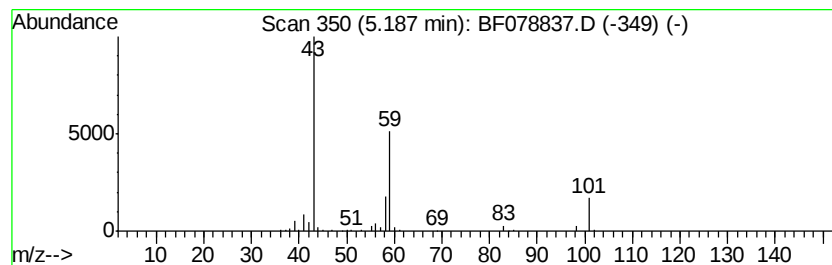
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.19	9.27 ng	220899	1,4-Dichlorobenzene-d4	7.38

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, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	25
3		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
4		5-Hexen-2-one	98	C6H10O	000109-49-9	9
5		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9



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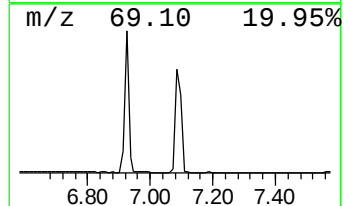
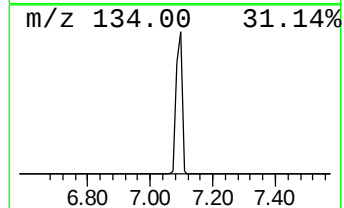
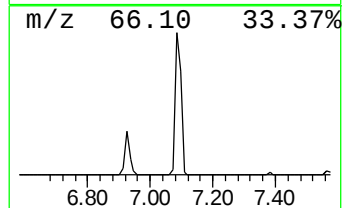
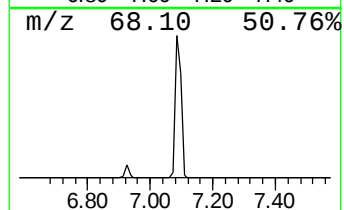
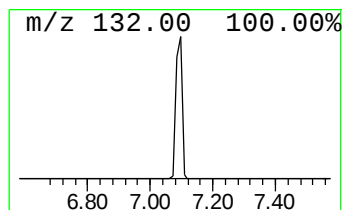
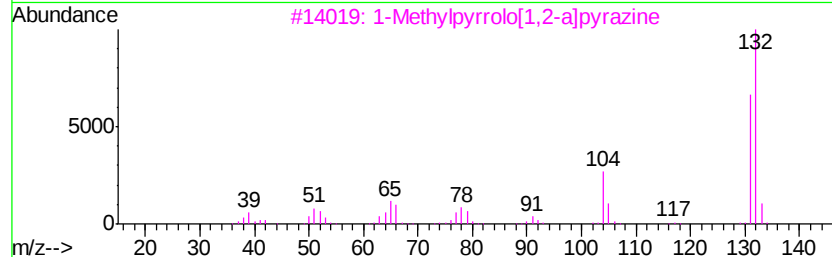
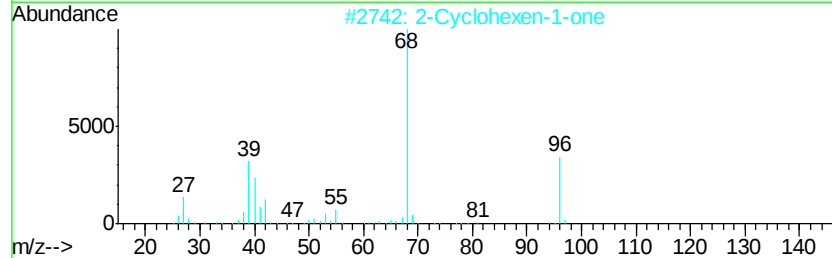
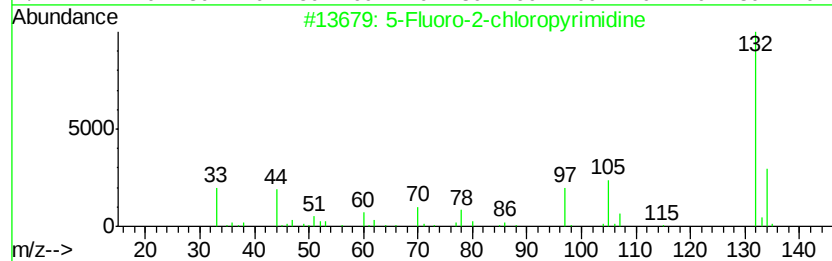
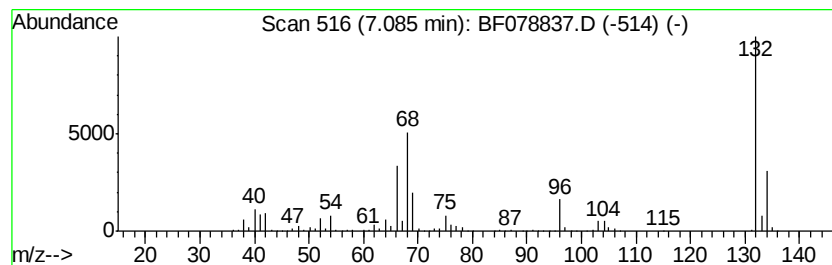
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 unknown7.08 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.08	64.21 ng	1530070	1,4-Dichlorobenzene-d4	7.38

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
2		2-Cyclohexen-1-one	96	C6H8O	000930-68-7	10
3		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9
4		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9
5		Xenon	132	Xe	007440-63-3	9



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042915\
Data File : BF078837.D
Acq On : 30 Apr 2015 4:41
Operator : TP/IZ
Sample : G2021-11
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
PS-25-0-2

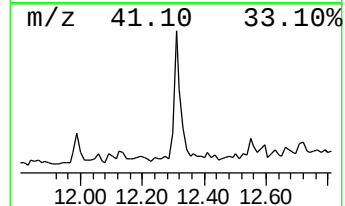
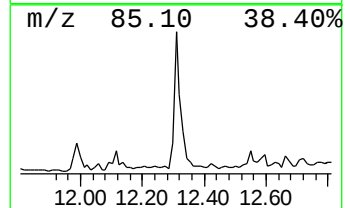
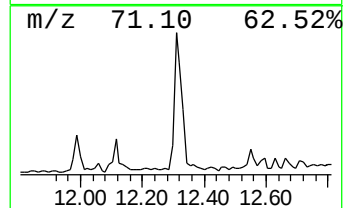
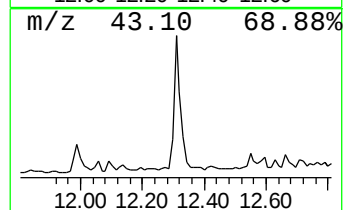
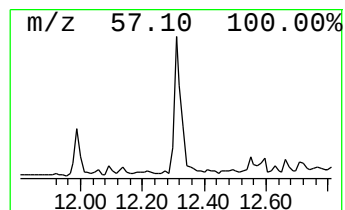
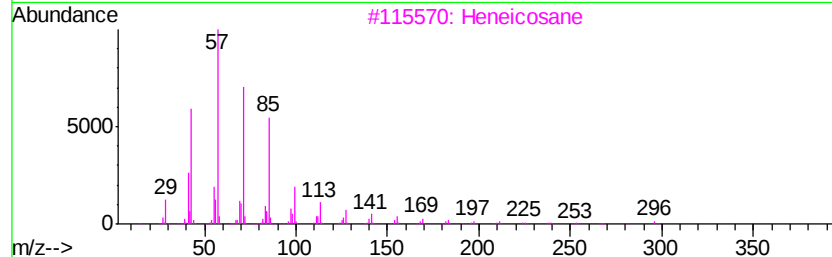
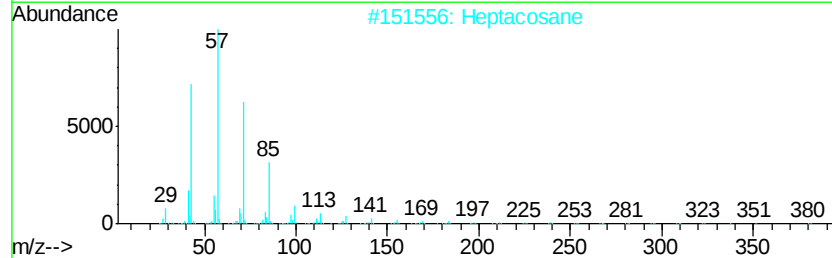
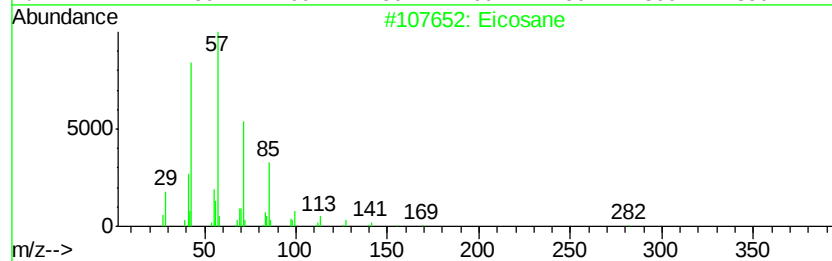
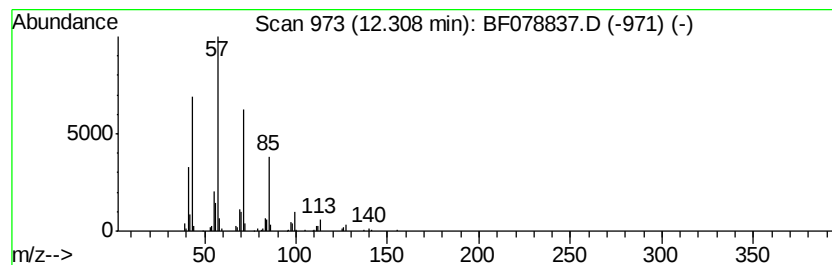
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Eicosane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.31	6.57 ng	297925	Phenanthrene-d10	12.98

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	91
2		Heptacosane	380	C27H56	000593-49-7	91
3		Heneicosane	296	C21H44	000629-94-7	90
4		Heptadecane	240	C17H36	000629-78-7	90
5		Hexadecane	226	C16H34	000544-76-3	90



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042915\
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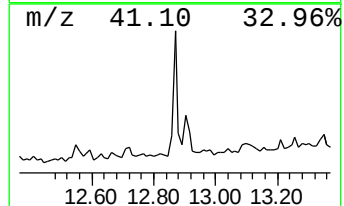
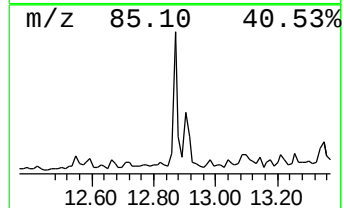
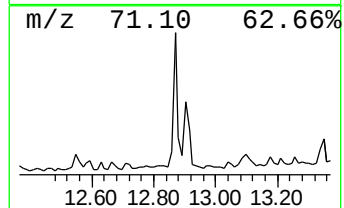
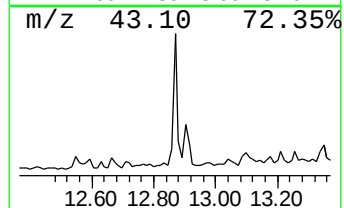
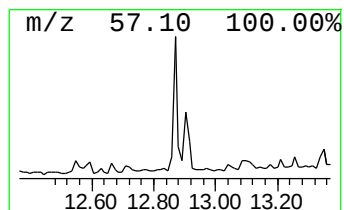
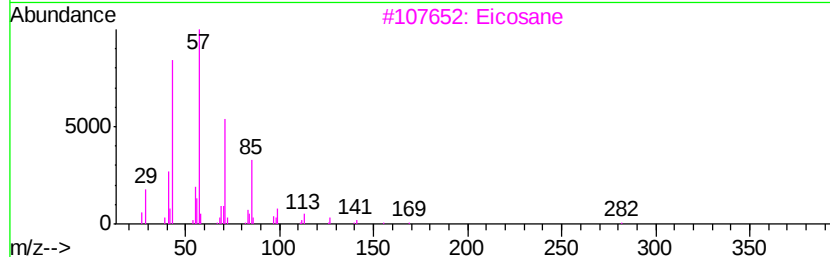
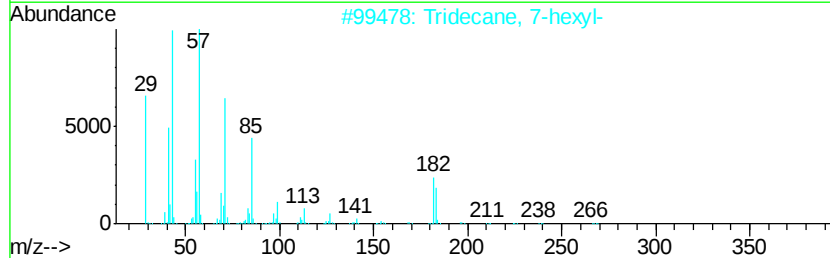
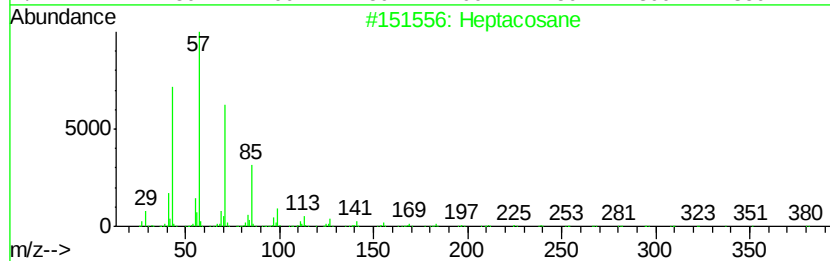
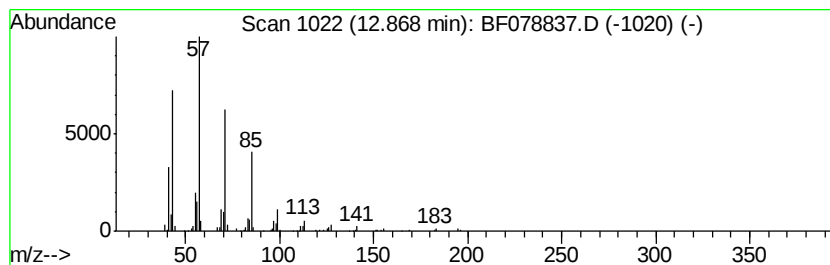
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Heptacosane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD		R.T.
12.87	5.20 ng	235660	Phenanthrene-d10		12.98
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptacosane	380	C27H56	000593-49-7	91
2	Tridecane, 7-hexyl-	268	C19H40	007225-66-3	91
3	Eicosane	282	C20H42	000112-95-8	91
4	Tridecane	184	C13H28	000629-50-5	90
5	Heneicosane	296	C21H44	000629-94-7	90



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042915\
Data File : BF078837.D
Acq On : 30 Apr 2015 4:41
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Misc :
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Instrument :
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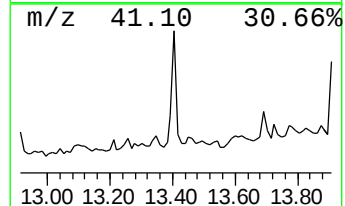
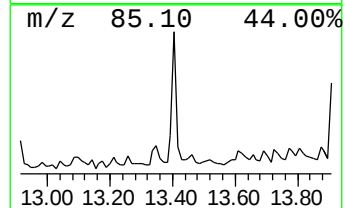
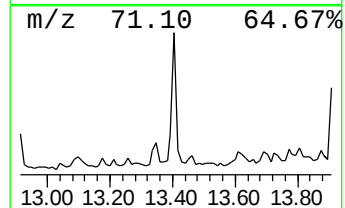
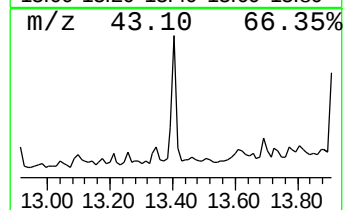
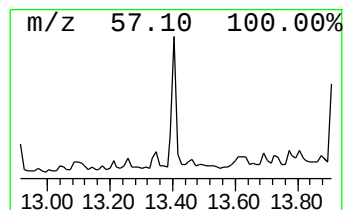
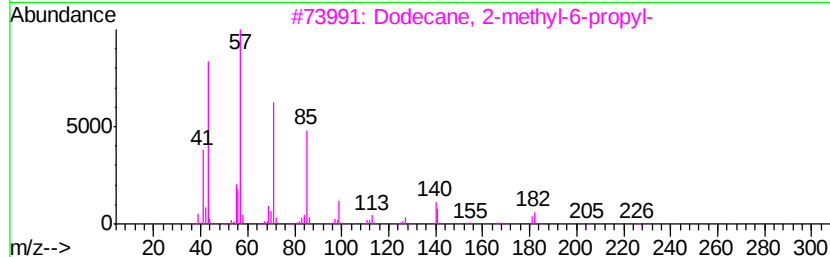
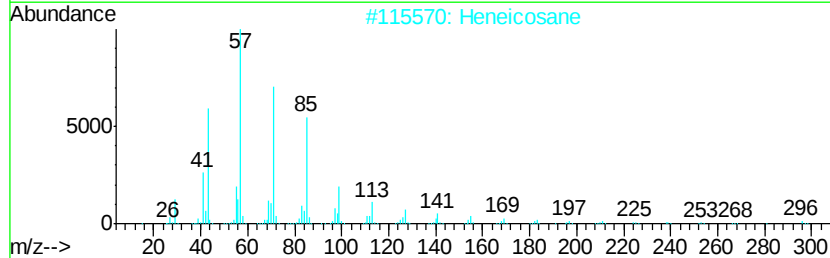
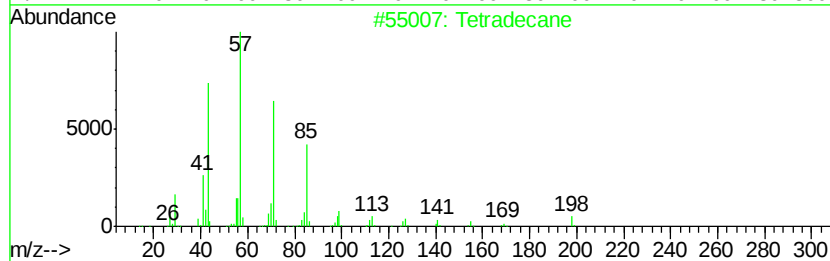
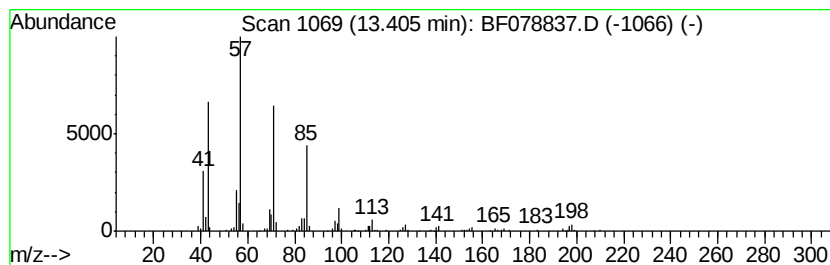
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Tetradecane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.41	5.18 ng	234736	Phenanthrene-d10	12.98

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetradecane	198	C14H30	000629-59-4	95
2			Heneicosane	296	C21H44	000629-94-7	90
3			Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	86
4			Nonadecane	268	C19H40	000629-92-5	83
5			Heptadecane, 8-methyl-	254	C18H38	013287-23-5	80



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042915\
Data File : BF078837.D
Acq On : 30 Apr 2015 4:41
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Misc :
ALS Vial : 31 Sample Multiplier: 1

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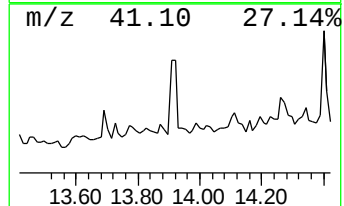
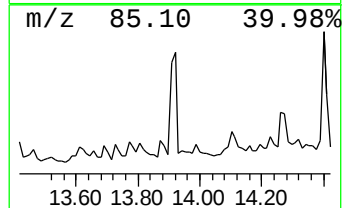
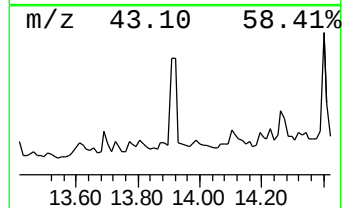
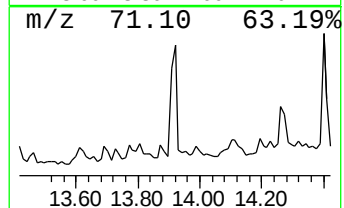
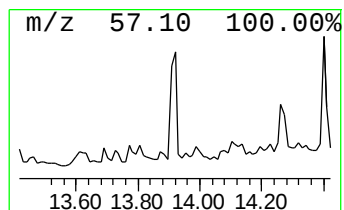
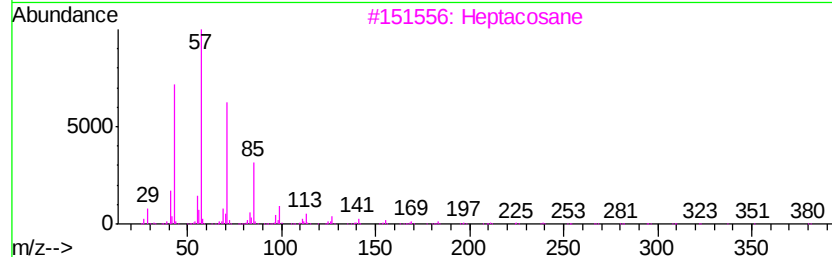
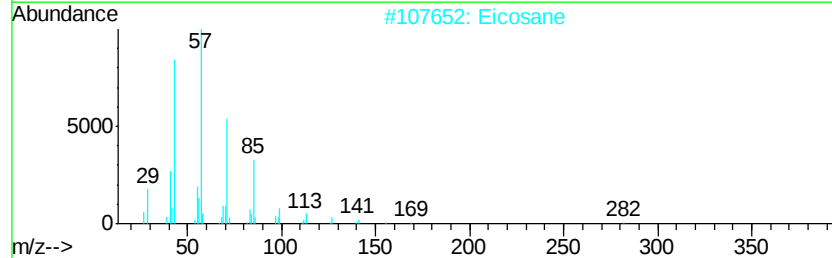
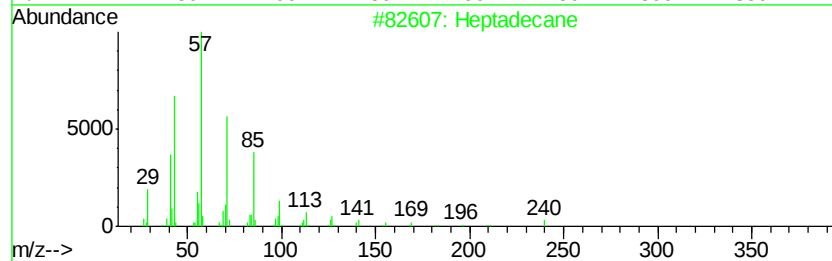
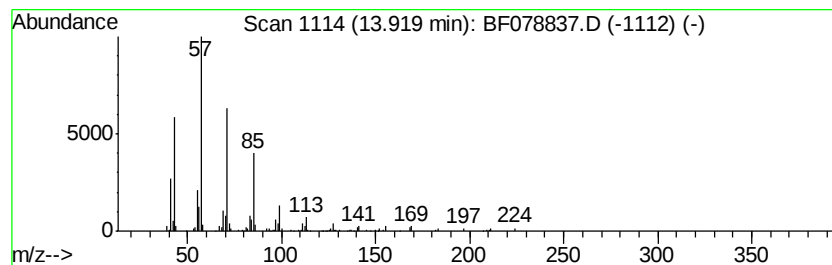
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Heptadecane Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.92	5.00 ng	226575	Phenanthrene-d10	12.98

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane	240	C17H36	000629-78-7	91
2		Eicosane	282	C20H42	000112-95-8	91
3		Heptacosane	380	C27H56	000593-49-7	91
4		Pentacosane	352	C25H52	000629-99-2	90
5		Heneicosane	296	C21H44	000629-94-7	90



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042915\
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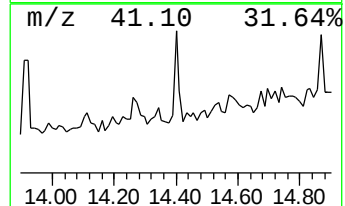
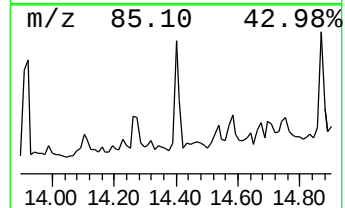
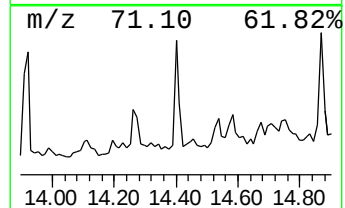
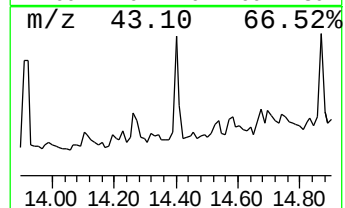
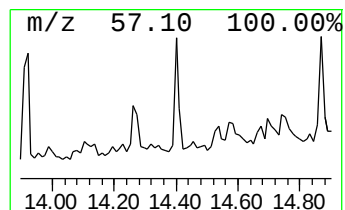
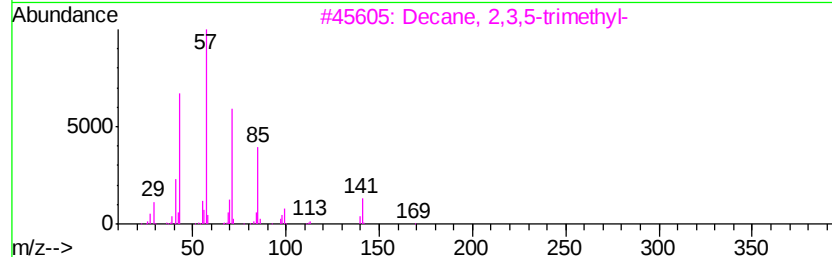
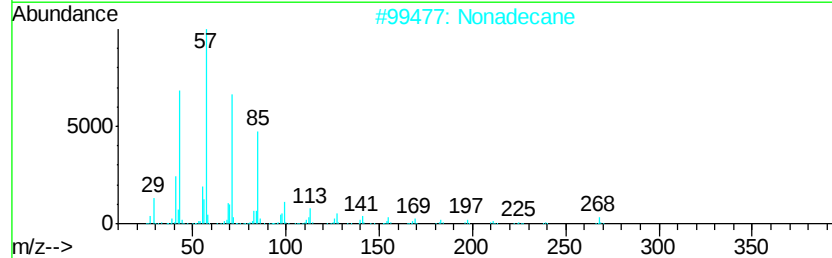
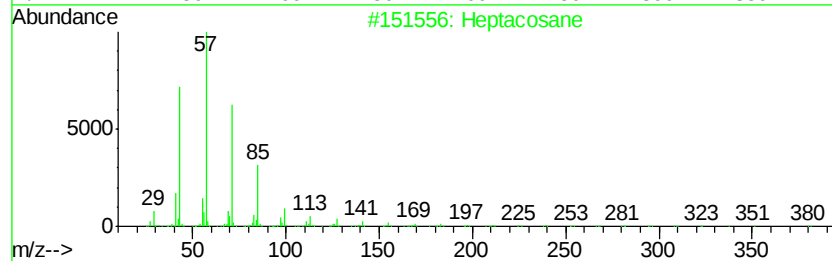
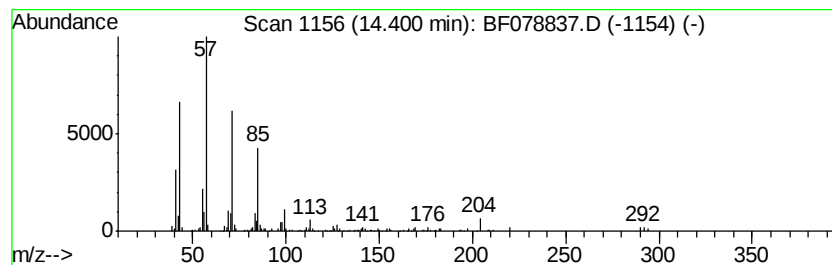
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Nonadecane Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.40	4.35 ng	197033	Phenanthrene-d10	12.98

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptacosane	380	C27H56	000593-49-7	83
2			Nonadecane	268	C19H40	000629-92-5	80
3			Decane, 2,3,5-trimethyl-	184	C13H28	062238-11-3	80
4			Pentadecane	212	C15H32	000629-62-9	80
5			Hexadecane	226	C16H34	000544-76-3	80



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042915\
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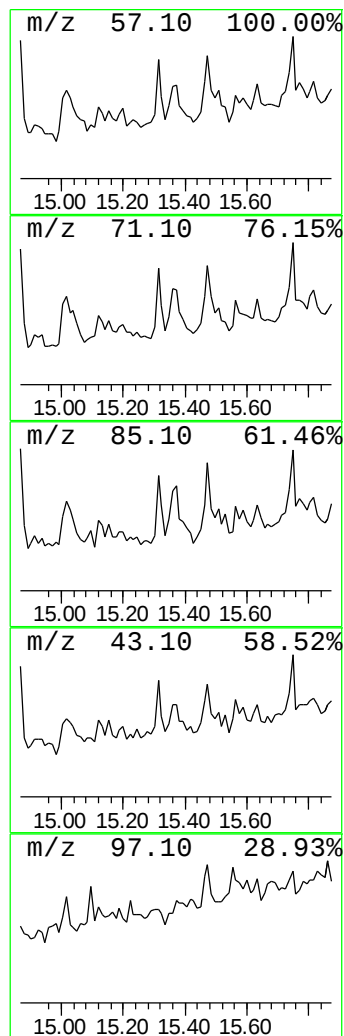
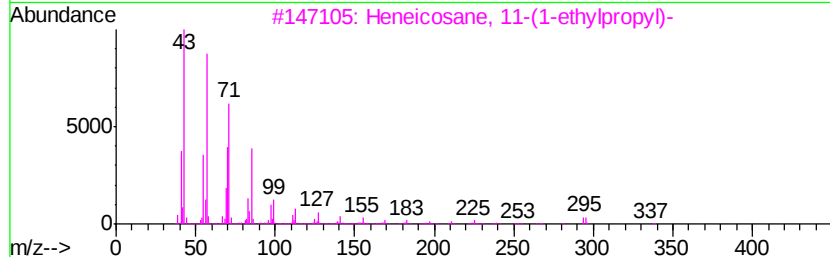
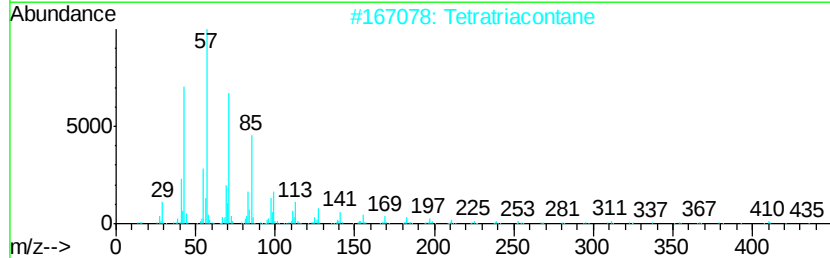
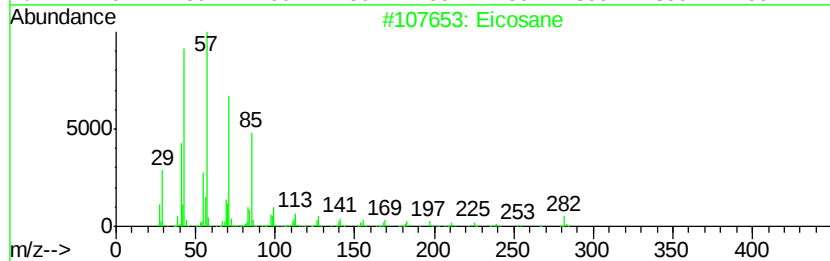
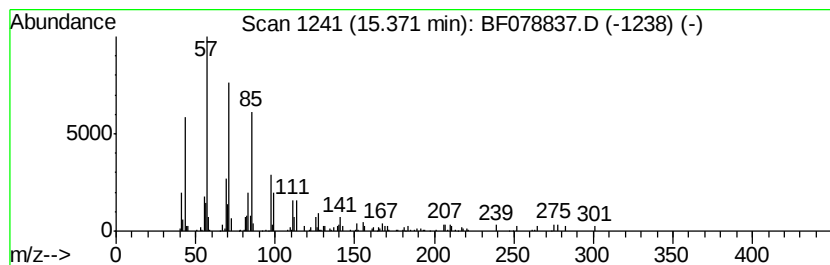
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 Tetratriacontane Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.37	2.68 ng	173060	Chrysene-d12	16.29
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Eicosane	282 C20H42	000112-95-8	89
2	Tetratriacontane	479 C34H70	014167-59-0	80
3	Heneicosane, 11-(1-ethylpropyl)-	366 C26H54	055282-11-6	80
4	Heneicosane	296 C21H44	000629-94-7	74
5	Heptadecane, 9-octyl-	352 C25H52	007225-64-1	72



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042915\
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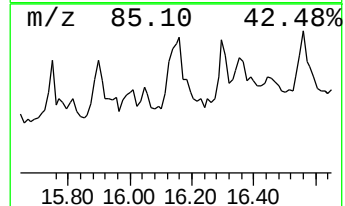
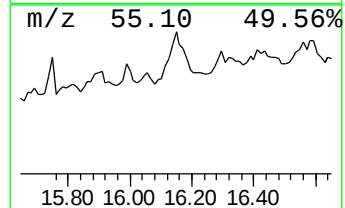
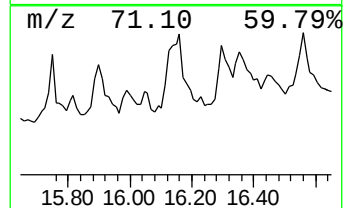
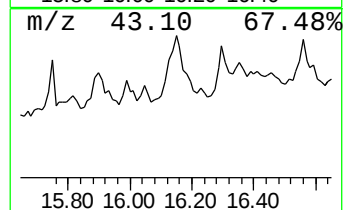
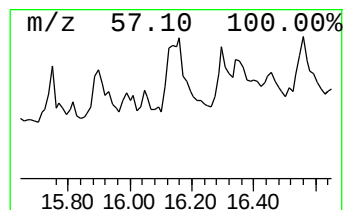
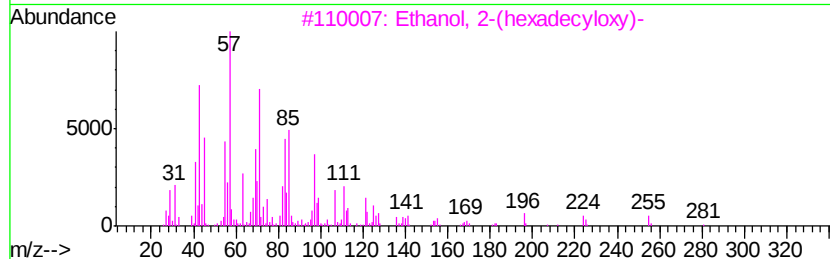
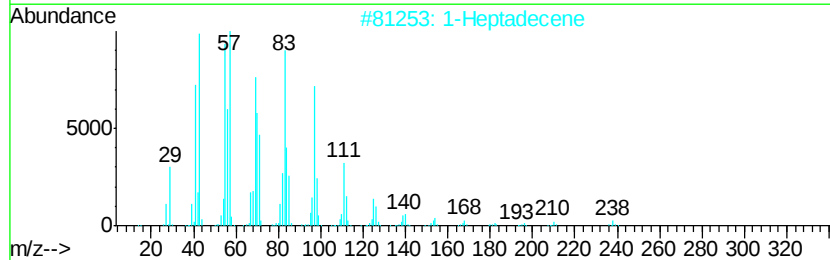
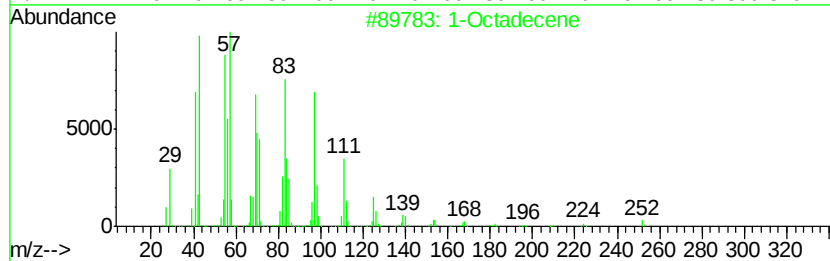
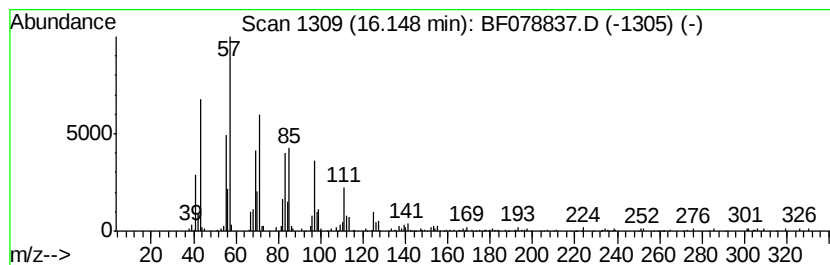
Instrument :
BNA_F
ClientSampleId :
PS-25-0-2

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF042815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 1-Octadecene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.15	11.30 ng	729082	Chrysene-d12	16.29
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1-Octadecene	252 C18H36	000112-88-9	92
2	1-Heptadecene	238 C17H34	006765-39-5	91
3	Ethanol, 2-(hexadecyloxy)-	286 C18H38O2	002136-71-2	91
4	Ethanol, 2-(octadecyloxy)-	314 C20H42O2	002136-72-3	74
5	Nonadecane, 1-chloro-	302 C19H39Cl	062016-76-6	70



Data Path : Z:\HPCHEM1\BNA F\DATA\BF042915\
Data File : BF078837.D
Acq On : 30 Apr 2015 4:41
Operator : TP/IZ
Sample : G2021-11
Misc :
ALS Vial : 31 Sample Multiplier: 1

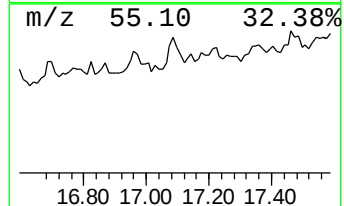
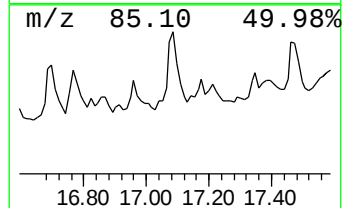
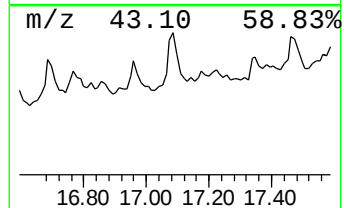
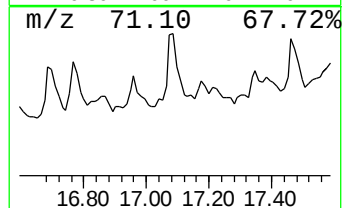
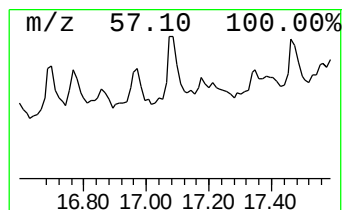
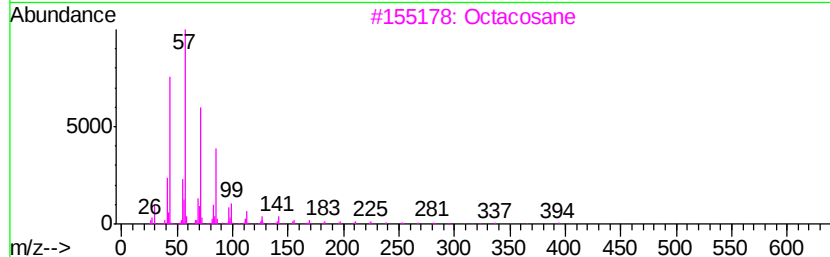
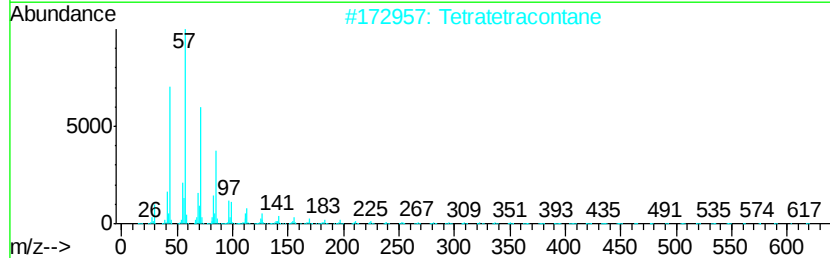
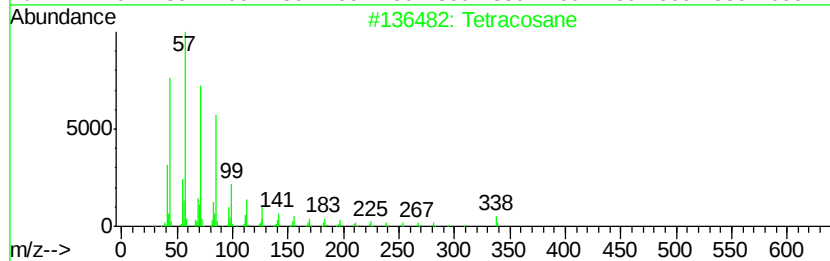
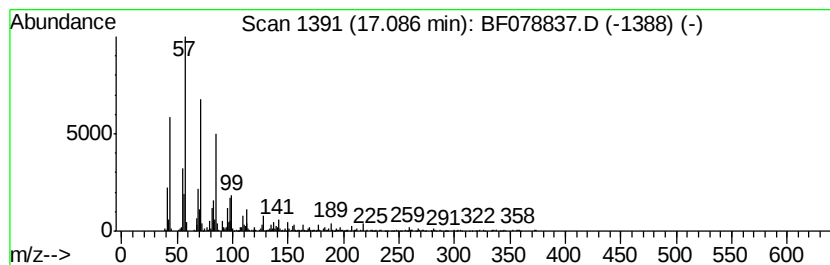
Instrument :
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ClientSampleId :
PS-25-0-2

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF042815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 Tetratetracontane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.09	8.09 ng	522031	Chrysene-d12	16.29
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Tetracosane	338 C24H50	000646-31-1 90
2		Tetratetracontane	619 C44H90	007098-22-8 83
3		Octacosane	394 C28H58	000630-02-4 83
4		Heptacosane	380 C27H56	000593-49-7 83
5		Tricosane, 2-methyl-	338 C24H50	001928-30-9 81



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042915\
Data File : BF078837.D
Acq On : 30 Apr 2015 4:41
Operator : TP/IZ
Sample : G2021-11
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
PS-25-0-2

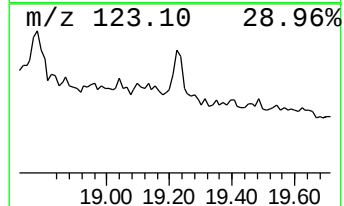
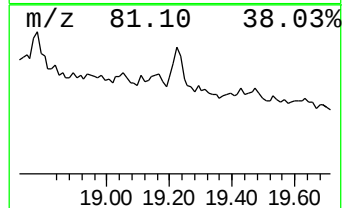
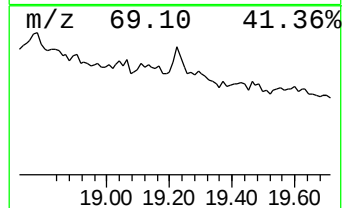
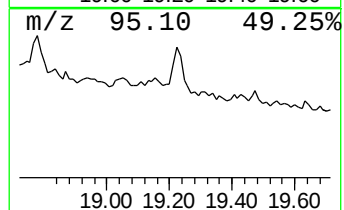
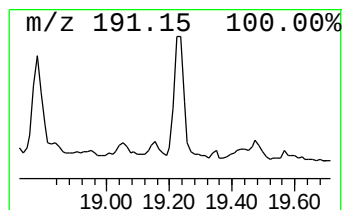
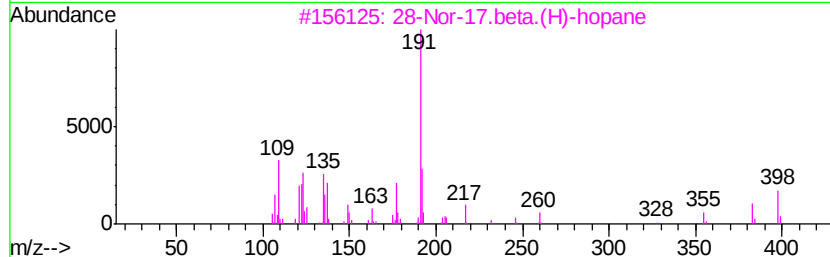
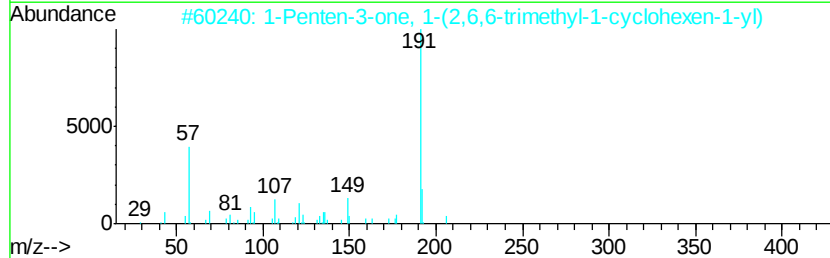
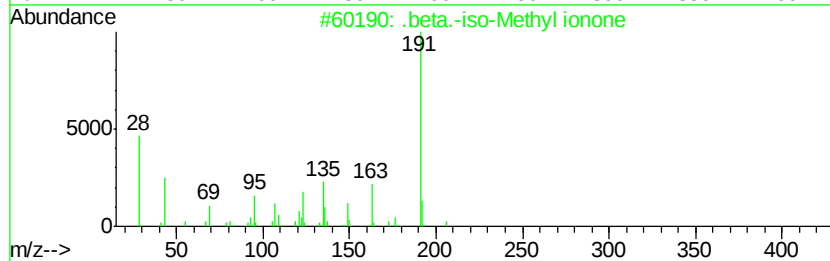
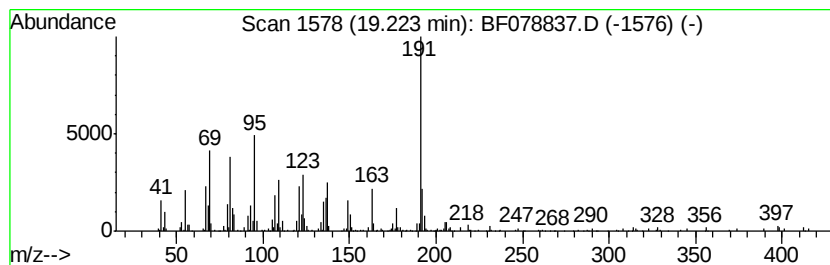
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 20 .beta.-iso-Methyl ionone Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.22	3.49 ng	267149	Perylene-d12	18.01

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	.beta.-iso-Methyl ionone	206	C14H22O	1000285-40-2	72
2		1-Penten-3-one, 1-(2,6,6-trimeth...	206	C14H22O	000127-43-5	43
3		28-Nor-17.beta.(H)-hopane	398	C29H50	036728-72-0	38
4		1-Cyclohexene, 1,3,3-trimethyl-2...	206	C14H22O	1000197-08-4	38
5		Anthracene, 9-dodecyltetradecahy...	360	C26H48	055401-75-7	38



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042915\
 Data File : BF078837.D
 Acq On : 30 Apr 2015 4:41
 Operator : TP/IZ
 Sample : G2021-11
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PS-25-0-2

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF042815.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
Cyclopentane, met...	1.29	108.7	ng	2589760	1	7.38	476589	20.0
Propane, 2,2-dime...	1.53	208.6	ng	4970050	1	7.38	476589	20.0
Butane, 2-methoxy...	1.86	24.4	ng	580727	1	7.38	476589	20.0
2-Pentanone, 4-hy...	5.19	9.3	ng	220899	1	7.38	476589	20.0
unknown7.08	7.08	64.2	ng	1530070	1	7.38	476589	20.0
Eicosane	12.31	6.6	ng	297925	4	12.98	906466	20.0
Heptacosane	12.87	5.2	ng	235660	4	12.98	906466	20.0
Tetradecane	13.41	5.2	ng	234736	4	12.98	906466	20.0
Heptadecane	13.92	5.0	ng	226575	4	12.98	906466	20.0
Nonadecane	14.40	4.3	ng	197033	4	12.98	906466	20.0
Tetratriacontane	15.37	2.7	ng	173060	5	16.29	1289880	20.0
1-Octadecene	16.15	11.3	ng	729082	5	16.29	1289880	20.0
Tetratetracontane	17.09	8.1	ng	522031	5	16.29	1289880	20.0
.beta.-iso-Methyl...	19.22	3.5	ng	267149	6	18.01	1531250	20.0