

Data Path : Z:\HPCHEM1\BNA F\DATA\BF031117\
 Data File : BF093581.D
 Acq On : 11 Mar 2017 21:38
 Operator : SJ/MA
 Sample : I2228-02 2X
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TEC-TP3

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-BF030717.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	4.858	258	260	273	rBV	318312	493015	16.57%	1.385%
2	5.316	298	300	316	rBV	1934223	2602219	87.48%	7.311%
3	6.378	391	393	401	rBV	2361857	2974745	100.00%	8.358%
4	6.493	401	403	406	rVB	2577885	2515049	84.55%	7.066%
5	6.721	420	423	425	rBV	1269767	1677399	56.39%	4.713%
6	6.870	434	436	441	rBV	1995151	1866576	62.75%	5.244%
7	6.973	444	445	448	rVB3	28154	44038	1.48%	0.124%
8	7.099	455	456	458	rBV2	21717	31033	1.04%	0.087%
9	7.259	465	470	471	rBV3	38990	72860	2.45%	0.205%
10	7.293	471	473	477	rBV	1816490	1708850	57.45%	4.801%
11	7.499	487	491	492	rBV4	36348	89712	3.02%	0.252%
12	7.521	492	493	495	rVB	70376	55998	1.88%	0.157%
13	7.579	495	498	499	rBV3	18637	35525	1.19%	0.100%
14	7.636	499	503	505	rVV3	66331	98995	3.33%	0.278%
15	7.681	505	507	509	rBV2	25592	37229	1.25%	0.105%
16	7.750	511	513	517	rVV4	30737	59451	2.00%	0.167%
17	7.819	517	519	522	rVV3	33079	52482	1.76%	0.147%
18	7.887	522	525	528	rBV3	28862	73562	2.47%	0.207%
19	8.002	533	535	538	rVV	1846248	2105145	70.77%	5.915%
20	8.070	540	541	542	rVB	147086	100901	3.39%	0.283%
21	8.093	542	543	547	rVB2	48595	82945	2.79%	0.233%
22	8.162	547	549	551	rBV3	46635	76071	2.56%	0.214%
23	8.207	551	553	554	rVB	95255	93218	3.13%	0.262%
24	8.287	558	560	564	rBV3	76898	138746	4.66%	0.390%
25	8.356	564	566	567	rBV	25592	35469	1.19%	0.100%
26	8.390	567	569	570	rVB	61855	58806	1.98%	0.165%
27	8.424	570	572	577	rBV2	212069	357805	12.03%	1.005%
28	8.504	577	579	581	rBV2	36554	77405	2.60%	0.217%
29	8.539	581	582	583	rBV	74847	54718	1.84%	0.154%
30	8.596	586	587	588	rBV	65353	70264	2.36%	0.197%
31	8.687	594	595	597	rVB2	128216	143421	4.82%	0.403%
32	8.733	597	599	600	rBV2	57915	107730	3.62%	0.303%
33	8.824	606	607	609	rBV2	178127	186494	6.27%	0.524%
34	8.950	616	618	620	rBV2	50704	117577	3.95%	0.330%

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Integration Parameters: rteint.p
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If leading or trailing edge < 100 prefer < Baseline drop else tangent >
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35	9.019	622	624	627	rBV2	201273	311868	10.48%	0.876%
36	9.076	627	629	631	rBV	1908342	1947173	65.46%	5.471%
37	9.144	633	635	638	rBV3	298219	451768	15.19%	1.269%
38	9.224	638	642	644	rVV3	102678	188134	6.32%	0.529%
39	9.350	651	653	654	rBV	92846	84791	2.85%	0.238%
40	9.453	660	662	663	rBV2	175647	209176	7.03%	0.588%
41	9.476	663	664	666	rVV2	529848	545952	18.35%	1.534%
42	9.567	670	672	674	rVB2	161642	187649	6.31%	0.527%
43	9.636	674	678	681	rBV5	343755	740893	24.91%	2.082%
44	9.693	681	683	684	rBV2	116965	182618	6.14%	0.513%
45	9.727	684	686	687	rBV	193868	276390	9.29%	0.777%
46	9.750	687	688	690	rVV	1525490	1886457	63.42%	5.300%
47	9.785	690	691	694	rVV3	79956	138770	4.66%	0.390%
48	10.013	709	711	712	rBV2	171961	221725	7.45%	0.623%
49	10.413	743	746	748	rVB2	266008	433094	14.56%	1.217%
50	10.550	756	758	760	rBV	885074	924943	31.09%	2.599%
51	10.676	766	769	772	rVB	546252	832340	27.98%	2.339%
52	11.133	808	809	813	rVB	437868	432623	14.54%	1.215%
53	11.236	816	818	822	rVV	1133290	1652161	55.54%	4.642%
54	12.825	954	957	959	rVB	1099891	1212830	40.77%	3.408%
55	13.854	1045	1047	1048	rBV	171027	234516	7.88%	0.659%
56	13.876	1048	1049	1052	rVB	1354737	1222978	41.11%	3.436%
57	14.391	1092	1094	1097	rBV	1065893	1185435	39.85%	3.331%
58	14.619	1112	1114	1117	rVB2	91913	139735	4.70%	0.393%
59	14.882	1135	1137	1139	rBV	110365	118170	3.97%	0.332%
60	15.271	1169	1171	1177	rVB	894761	1288561	43.32%	3.620%
61	15.648	1202	1204	1206	rVB	47643	60393	2.03%	0.170%
62	18.871	1482	1486	1491	rVB4	61263	185861	6.25%	0.522%

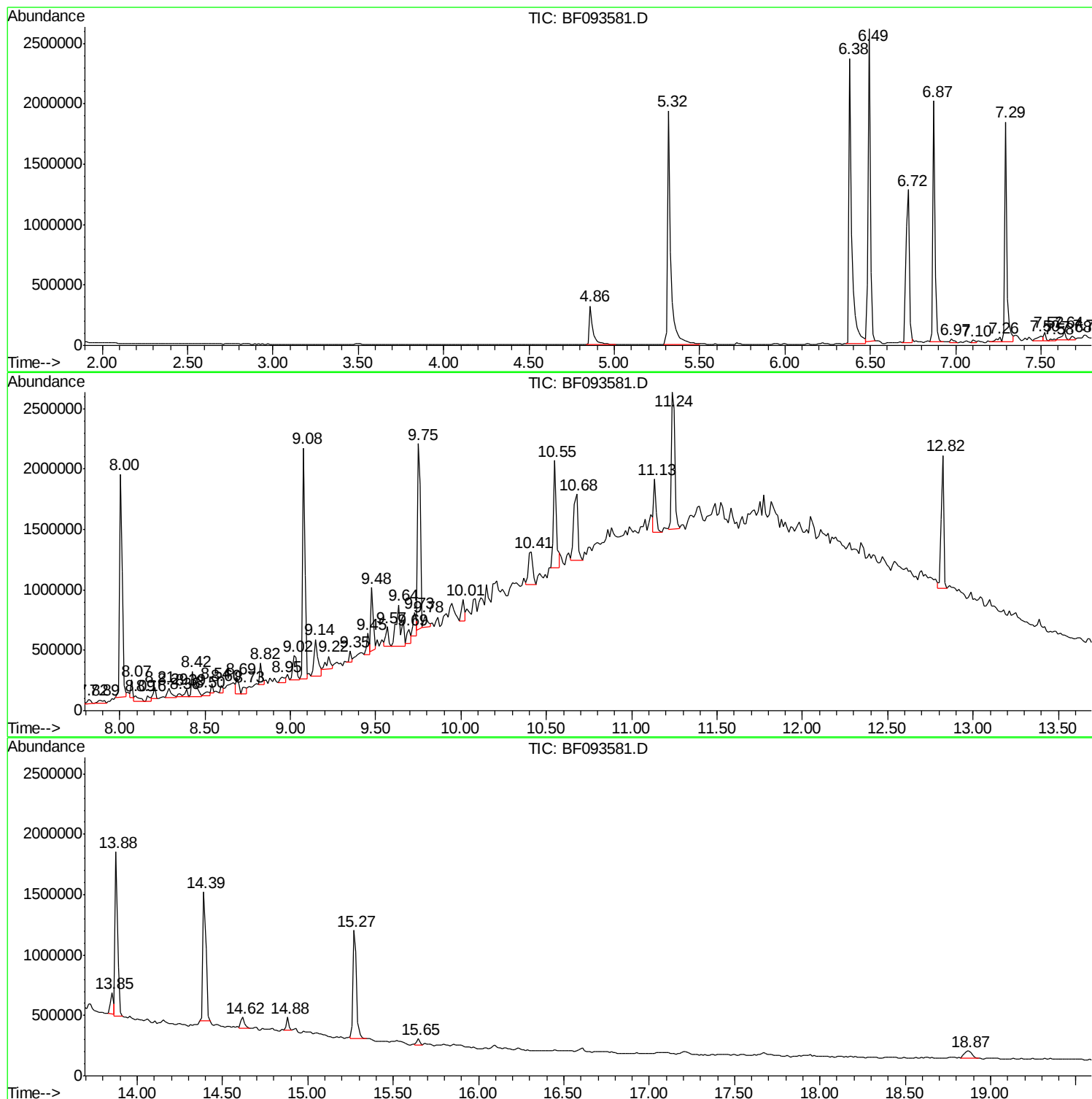
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF030717.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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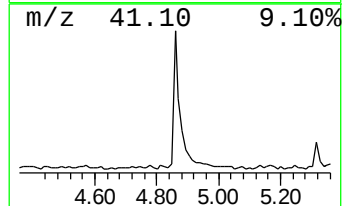
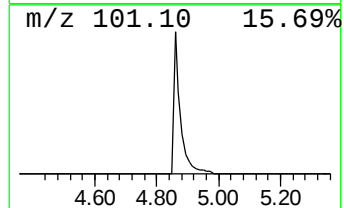
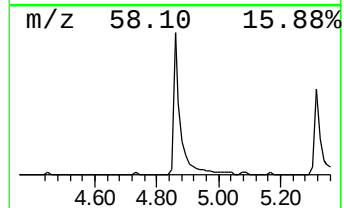
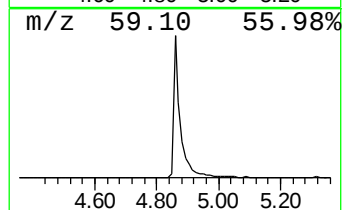
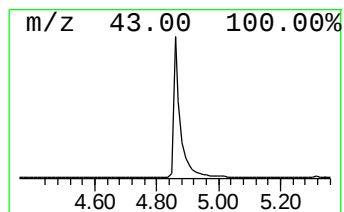
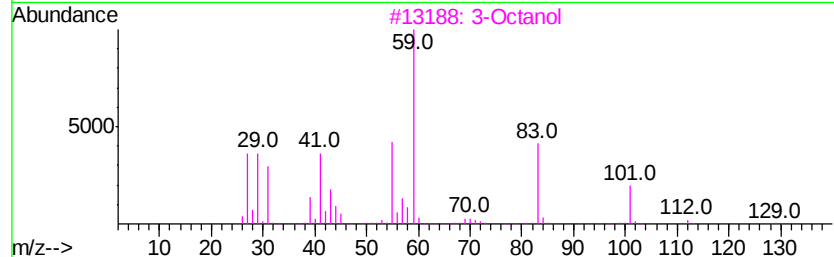
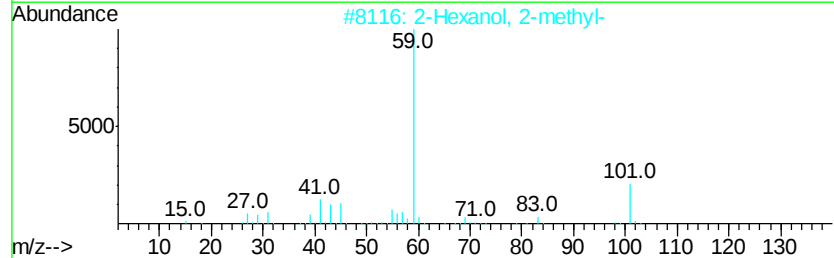
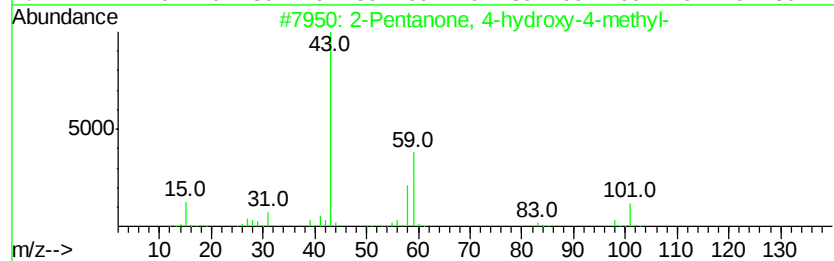
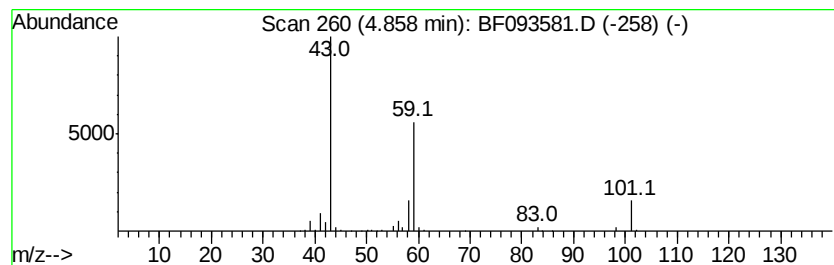
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF030717.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.86	5.88 ng	493015	1,4-Dichlorobenzene-d4	6.72

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	33
3			3-Octanol	130	C8H18O	000589-98-0	33
4			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
5			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	32



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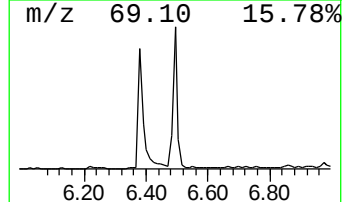
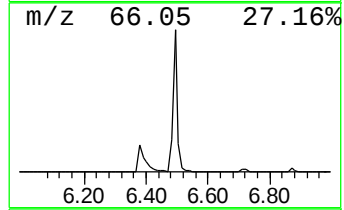
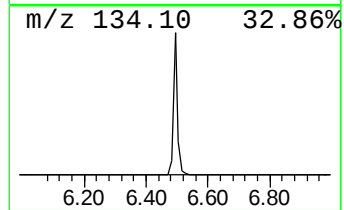
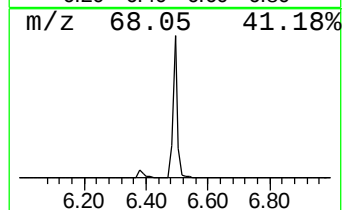
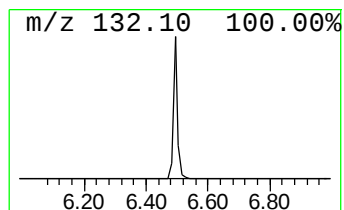
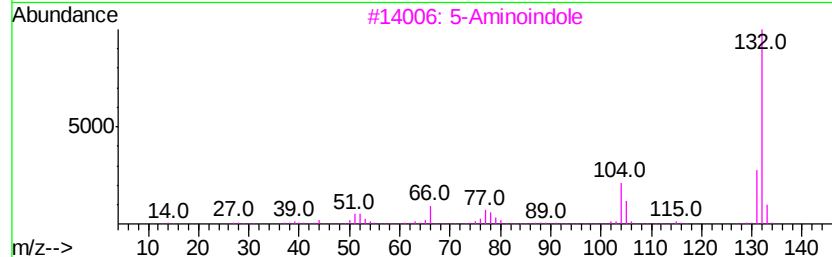
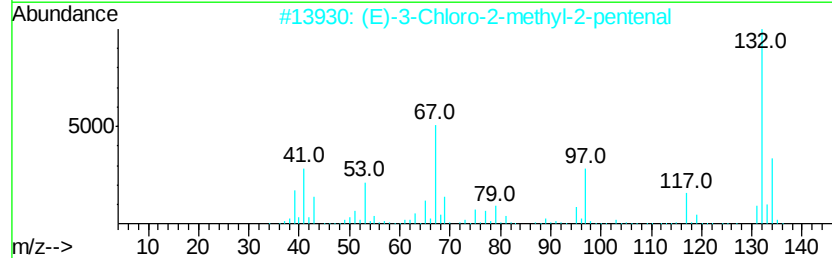
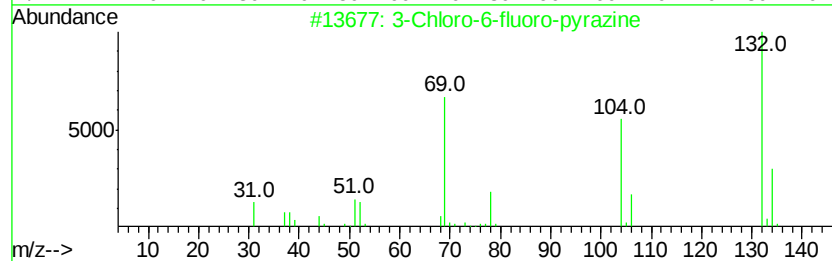
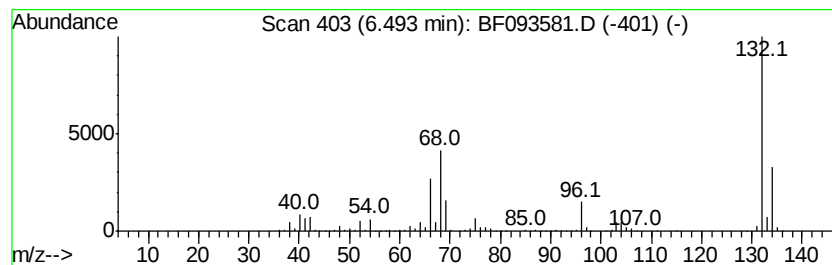
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown6.49 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.49	29.99 ng	2515050	1,4-Dichlorobenzene-d4	6.72

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	27
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9



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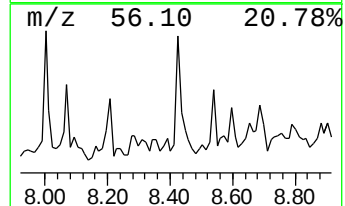
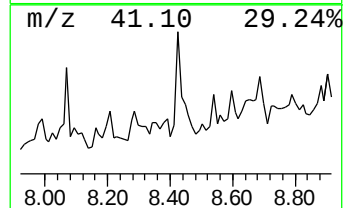
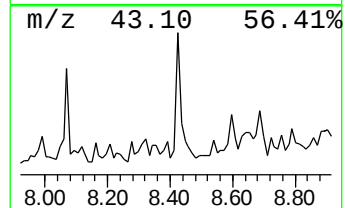
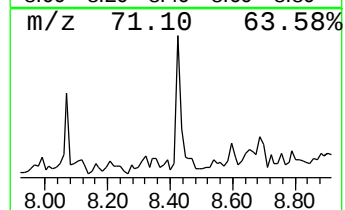
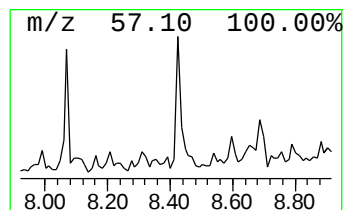
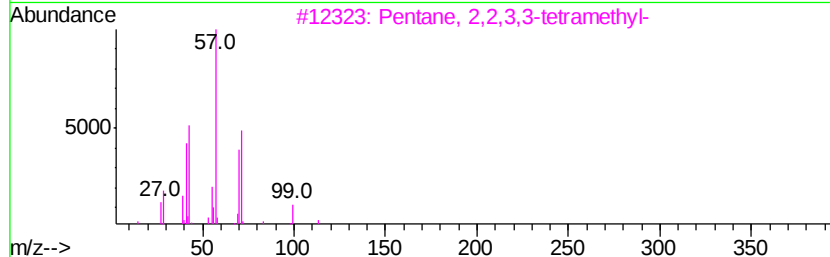
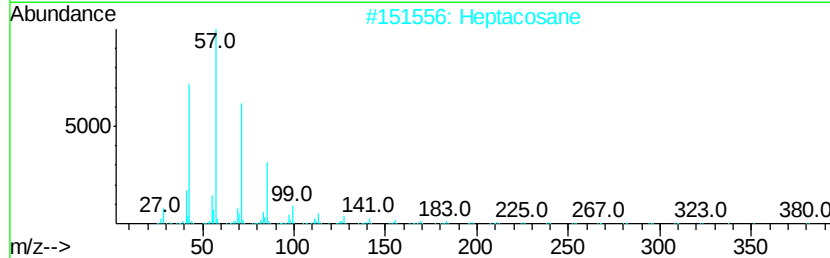
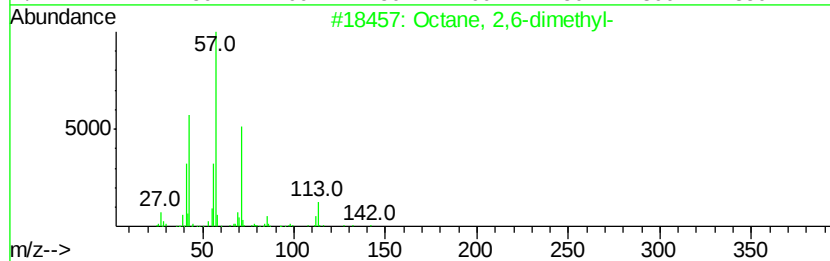
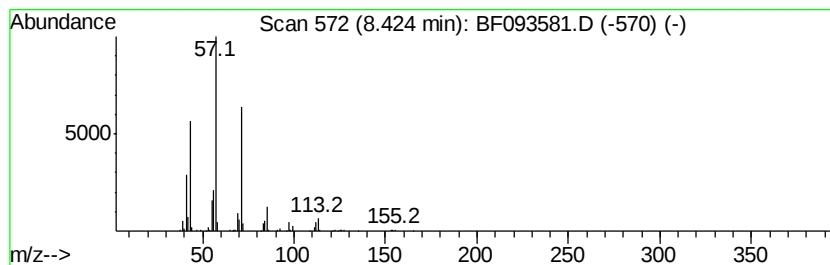
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 Octane, 2,6-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.42	3.40 ng	357805	Napthalene-d8	8.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	64
2		Heptacosane	380	C27H56	000593-49-7	53
3		Pentane, 2,2,3,3-tetramethyl-	128	C9H20	007154-79-2	53
4		Heptadecane, 2,6,10,14-tetramethyl-	296	C21H44	018344-37-1	53
5		Hexadecane	226	C16H34	000544-76-3	53



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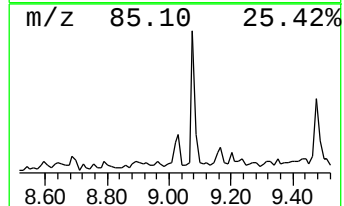
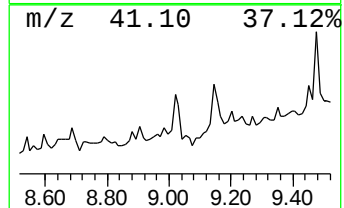
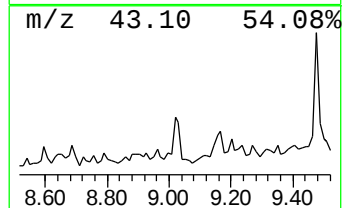
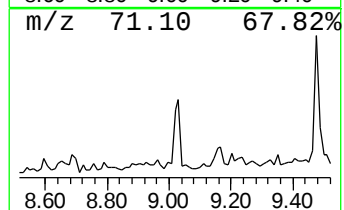
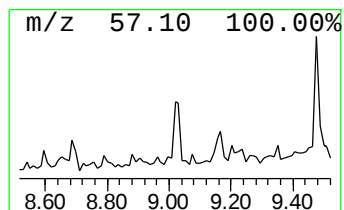
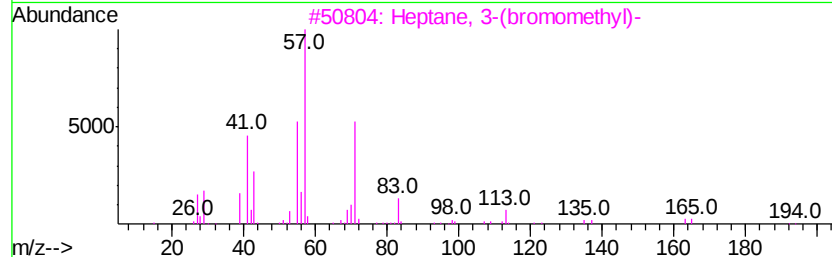
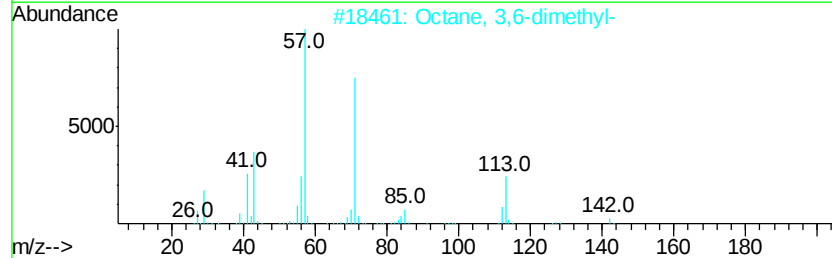
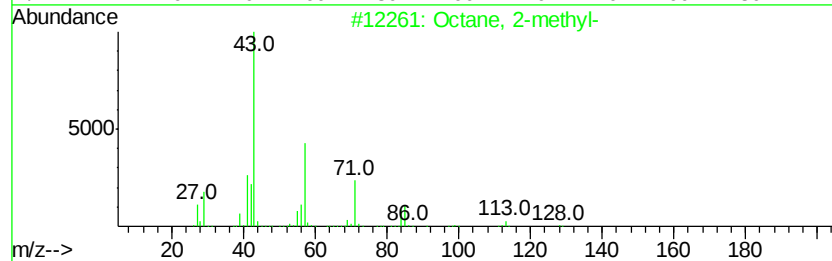
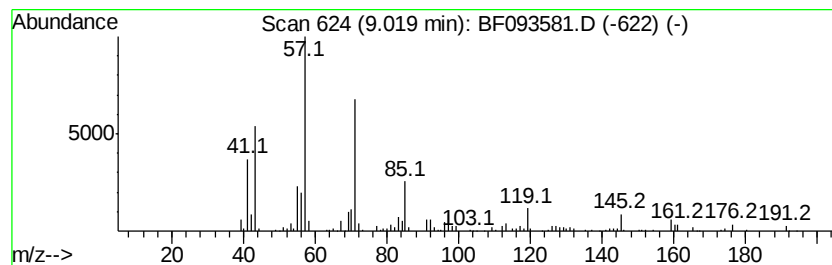
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Octane, 2-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.02	3.31 ng	311868	Acenaphthene-d10	9.75

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octane, 2-methyl-	128	C9H20	003221-61-2	59
2		Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	50
3		Heptane, 3-(bromomethyl)-	192	C8H17Br	018908-66-2	49
4		Undecane, 3,9-dimethyl-	184	C13H28	017301-31-4	47
5		Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	47



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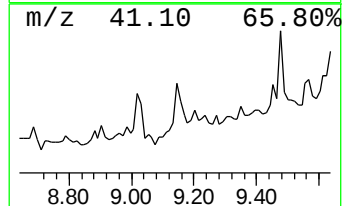
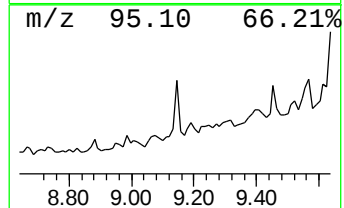
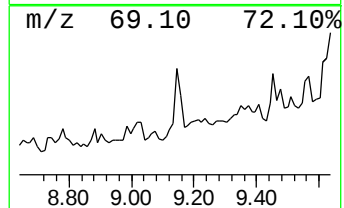
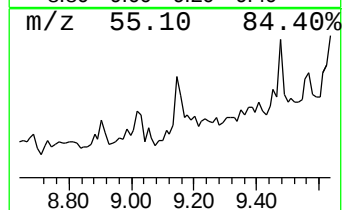
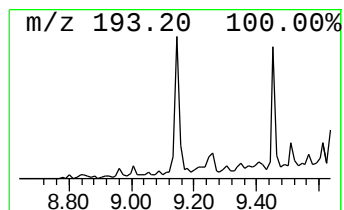
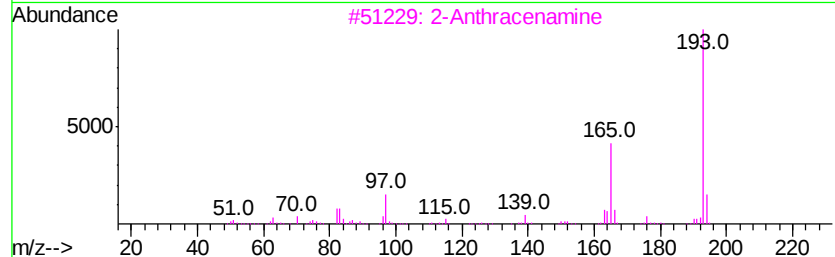
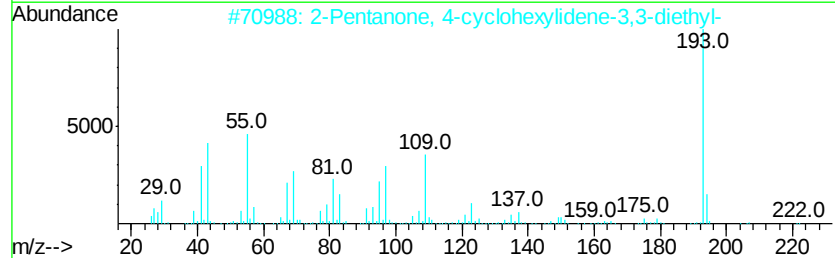
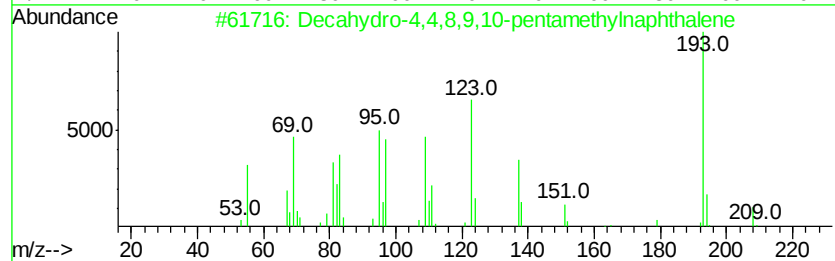
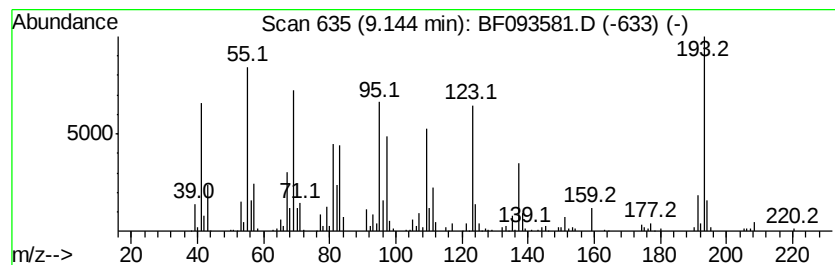
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Decahydro-4,4,8,9,10-pentam... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.14	4.79 ng	451768	Acenaphthene-d10	9.75

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Decahydro-4,4,8,9,10-pentamethyl...	208	C15H28	080655-44-3	90
2		2-Pentanone, 4-cyclohexylidene-3...	222	C15H26O	313253-65-5	38
3		2-Anthracenamine	193	C14H11N	000613-13-8	22
4		1,7-Dimethyl-4-(1-methylethyl)cy...	210	C15H30	000645-10-3	18
5		Naphthalene, 2-(1,1-dimethylethy...	208	C15H28	054934-96-2	18



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF031117\
Data File : BF093581.D
Acq On : 11 Mar 2017 21:38
Operator : SJ/MA
Sample : I2228-02 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

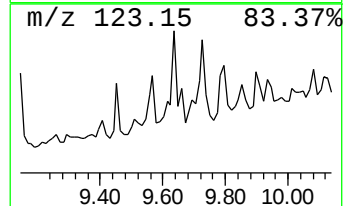
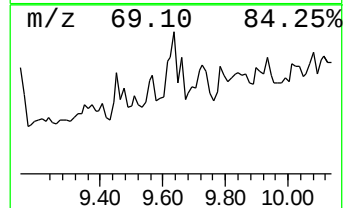
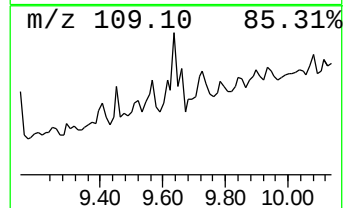
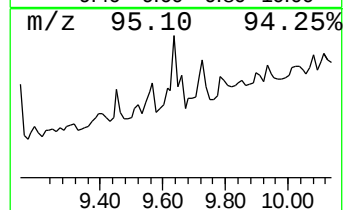
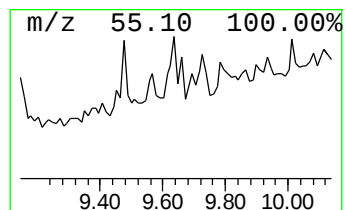
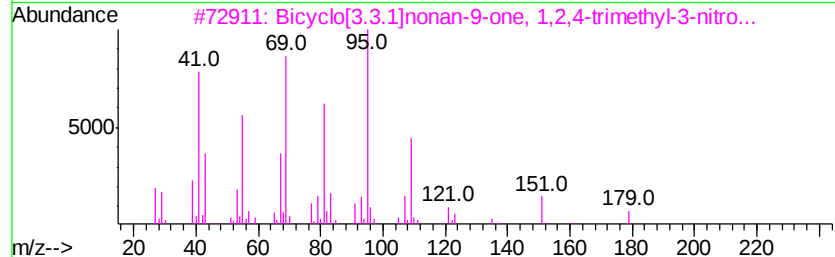
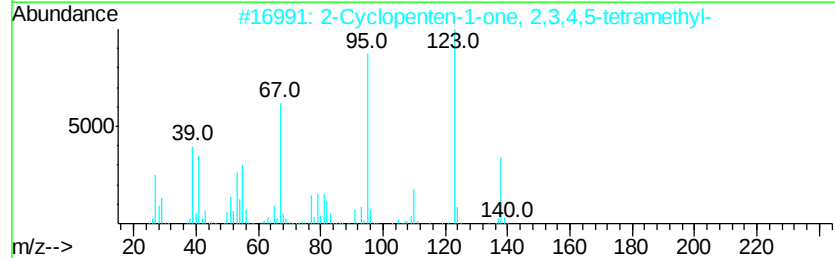
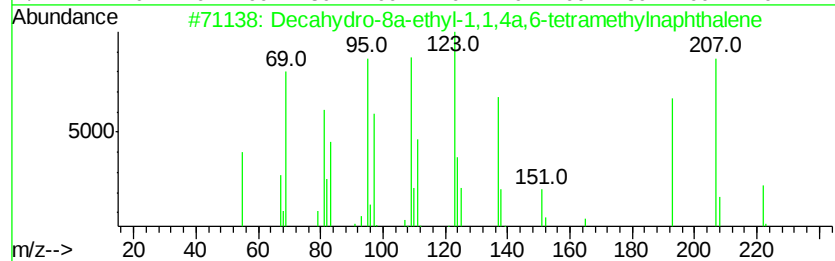
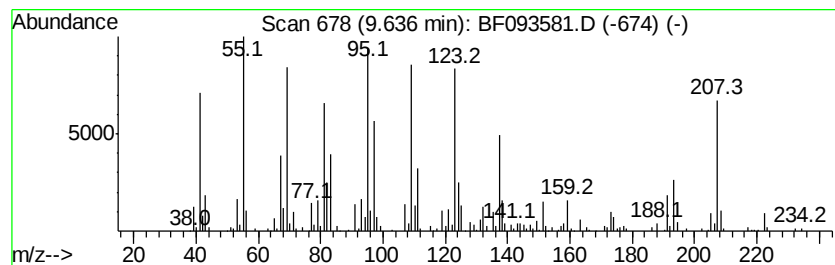
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF030717.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 unknown9.64 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
9.64	7.85 ng	740893	Acenaphthene-d10	9.75	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decahydro-8a-ethyl-1,1,4a,6-tetr...	222	C16H30	1000100-23-6	46
2	2-Cyclopenten-1-one, 2,3,4,5-tet...	138	C9H14O	054458-61-6	35
3	Bicyclo[3.3.1]nonan-9-one, 1,2,4...	225	C12H19NO3	129967-65-3	30
4	Naphthalene, decahydro-1,6-dimet...	208	C15H28	034315-85-0	27
5	Cyclopentene, 1-isopropyl-2,3-di...	138	C10H18	007712-73-4	25



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF031117\
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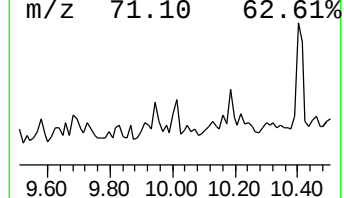
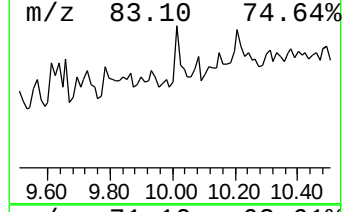
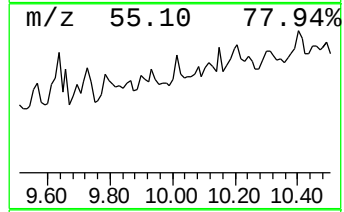
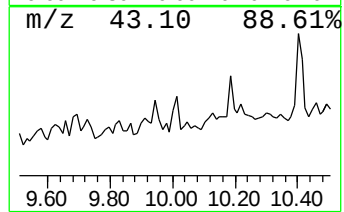
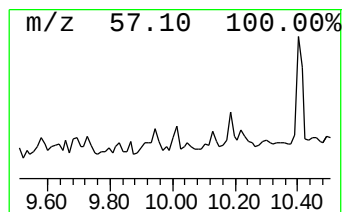
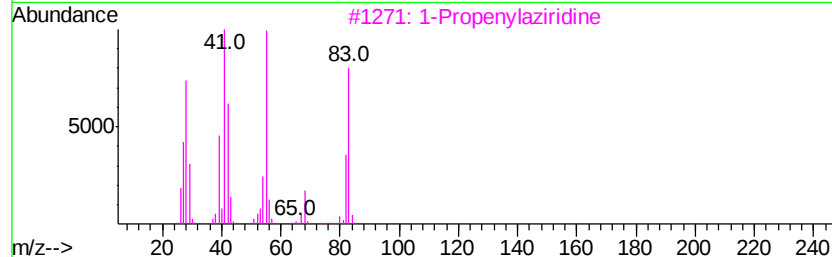
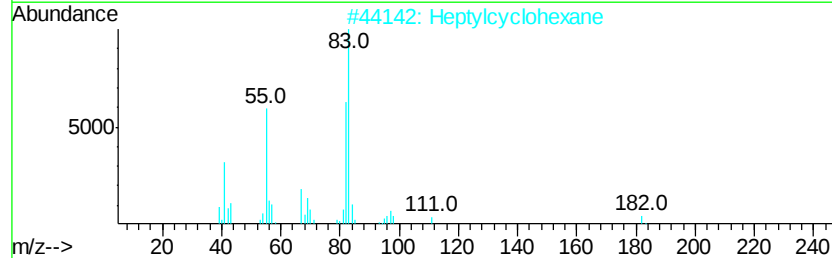
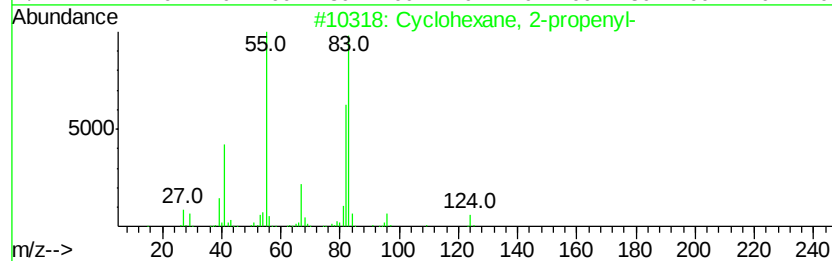
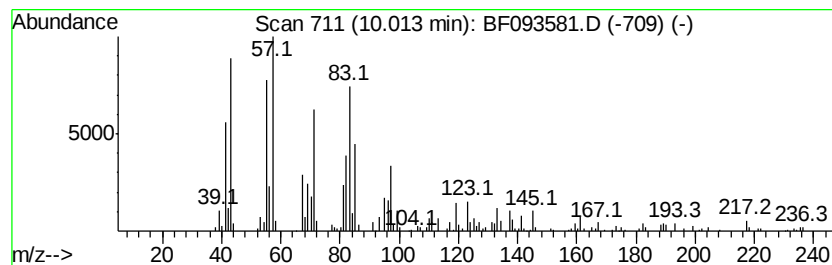
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF030717.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 unknown10.01 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD		R.T.
10.01	2.35 ng	221725	Acenaphthene-d10		9.75
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cvclohexane, 2-propenvl-	124	C9H16	002114-42-3	38
2	Heptylcvclohexane	182	C13H26	005617-41-4	38
3	1-Propenvlaziridine	83	C5H9N	1000192-51-0	38
4	2-Piperidinone, N-[4-bromo-n-but...	233	C9H16BrNO	195194-80-0	35
5	Bicyclo[3.1.1]heptan-2-one, 6,6-...	138	C9H14O	024903-95-5	25



Data Path : Z:\HPCHEM1\BNA F\DATA\BF031117\
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Sample : I2228-02 2X
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ALS Vial : 18 Sample Multiplier: 1

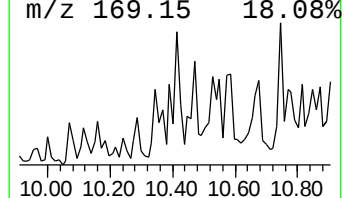
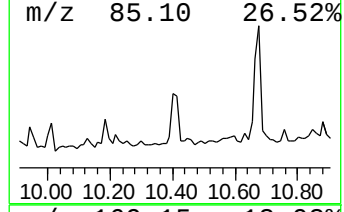
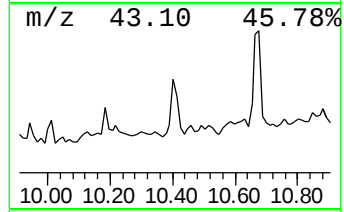
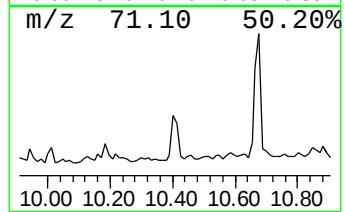
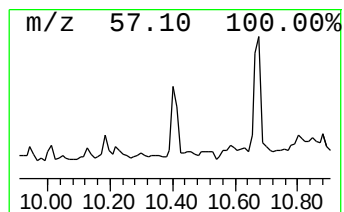
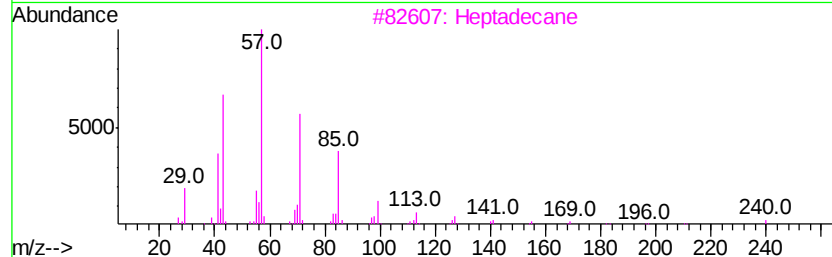
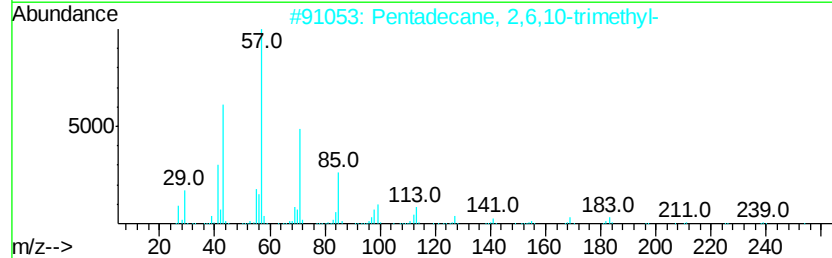
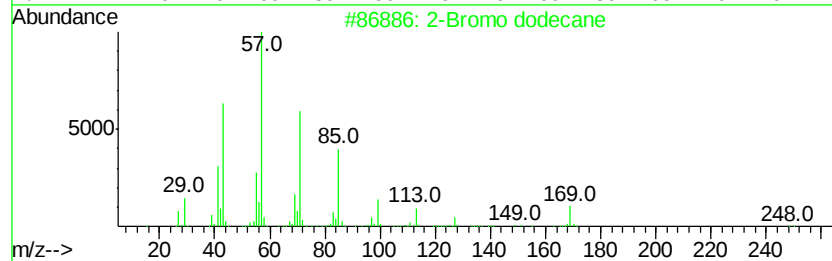
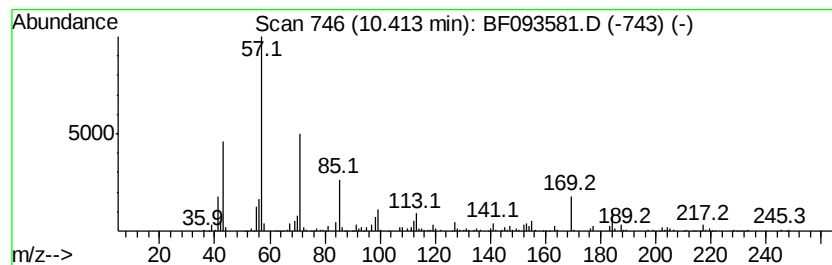
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF030717.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 2-Bromo dodecane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD		R.T.
10.41	4.59 ng	433094	Acenaphthene-d10		9.75
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Bromo dodecane	248	C12H25Br	013187-99-0	81
2	Pentadecane, 2,6,10-trimethyl-	254	C18H38	003892-00-0	64
3	Heptadecane	240	C17H36	000629-78-7	64
4	Hexadecane	226	C16H34	000544-76-3	59
5	Pentadecane	212	C15H32	000629-62-9	53



Data Path : Z:\HPCHEM1\BNA F\DATA\BF031117\
Data File : BF093581.D
Acq On : 11 Mar 2017 21:38
Operator : SJ/MA
Sample : I2228-02 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

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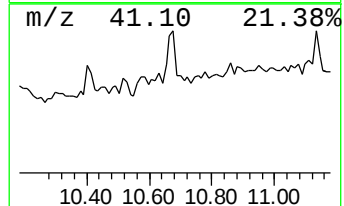
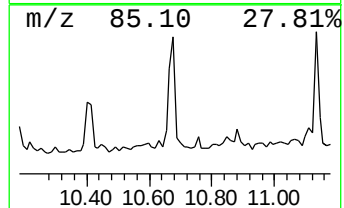
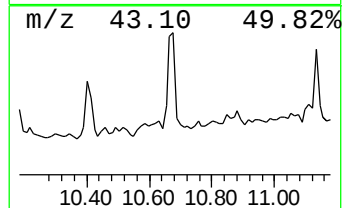
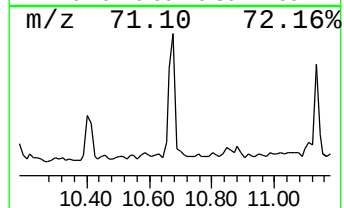
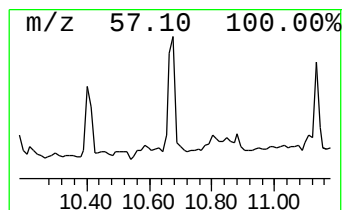
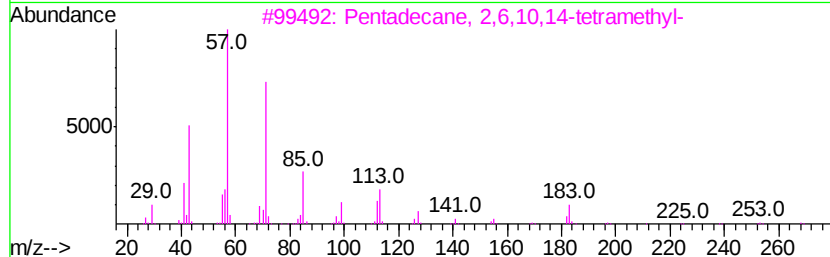
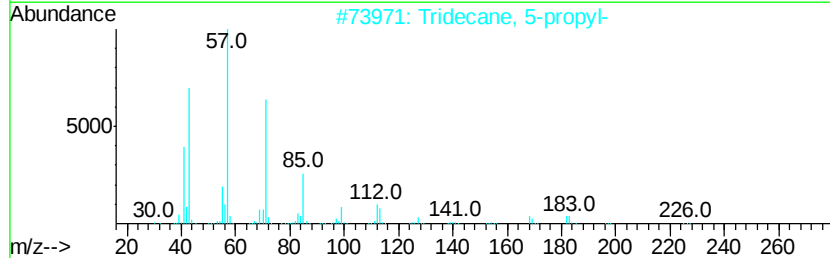
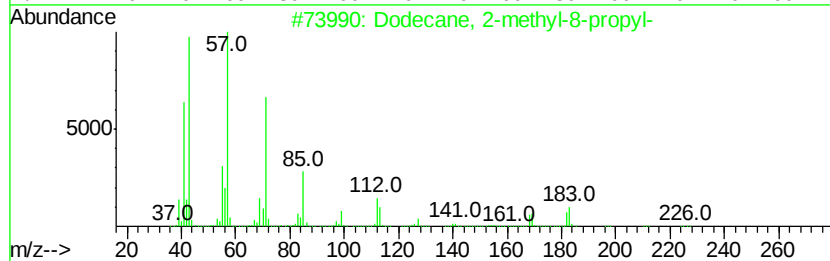
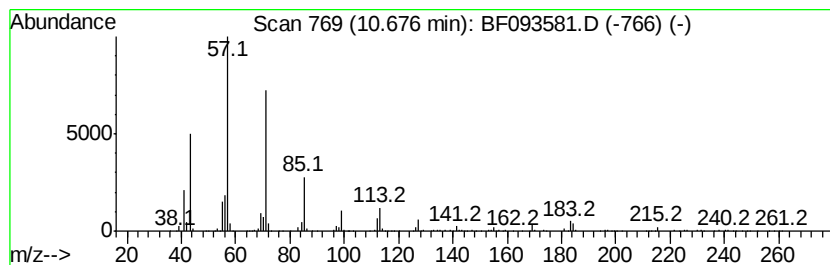
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Dodecane, 2-methyl-8-propyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.68	10.08 ng	832340	Phenanthrene-d10	11.24

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Dodecane, 2-methyl-8-propyl-	226	C16H34	055045-07-3	80
2		Tridecane, 5-propyl-	226	C16H34	055045-11-9	78
3		Pentadecane, 2,6,10,14-tetramethyl-	268	C19H40	001921-70-6	64
4		Dodecane, 2,6,10-trimethyl-	212	C15H32	003891-98-3	64
5		5,5-Dibutylnonane	240	C17H36	006008-17-9	64



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF031117\
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Operator : SJ/MA
Sample : I2228-02 2X
Misc :
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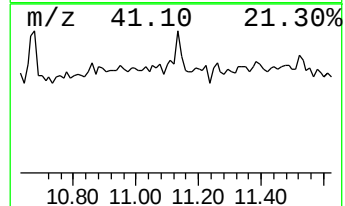
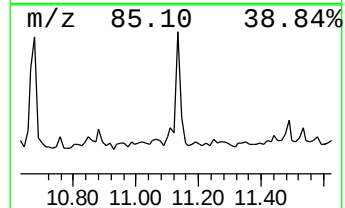
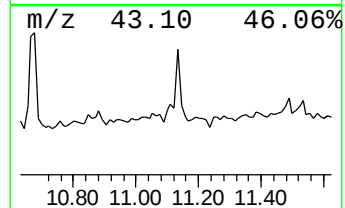
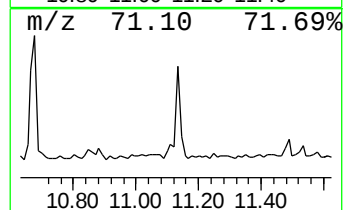
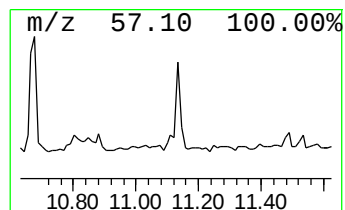
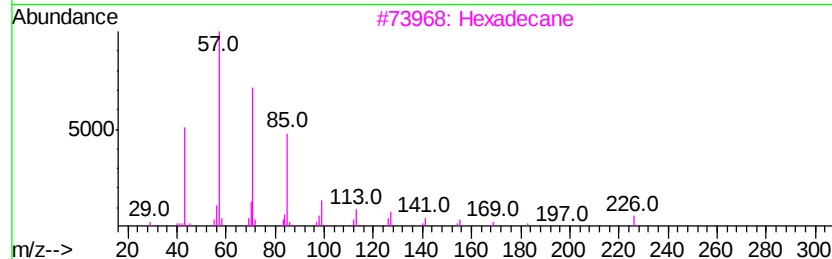
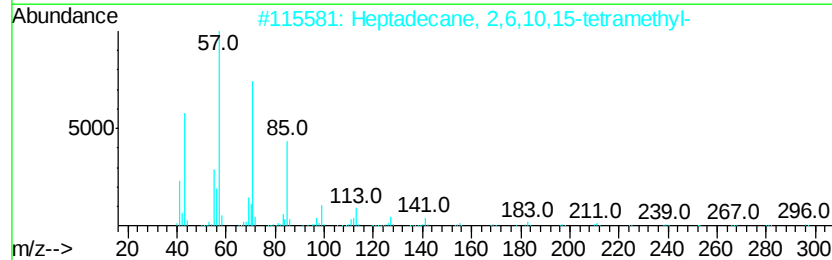
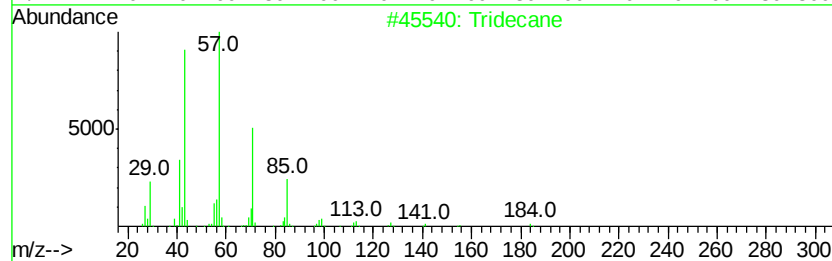
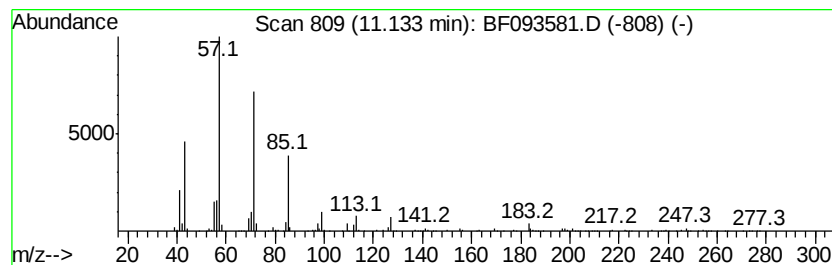
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF030717.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Tridecane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD		R.T.
11.13	5.24 ng	432623	Phenanthrene-d10		11.24
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecane	184	C13H28	000629-50-5	90
2	Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	86
3	Hexadecane	226	C16H34	000544-76-3	78
4	Heptacosane	380	C27H56	000593-49-7	78
5	Decane, 3-methyl-	156	C11H24	013151-34-3	72



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF031117\
Data File : BF093581.D
Acq On : 11 Mar 2017 21:38
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Misc :
ALS Vial : 18 Sample Multiplier: 1

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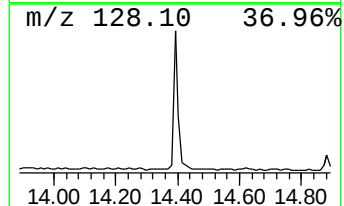
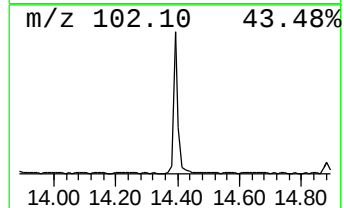
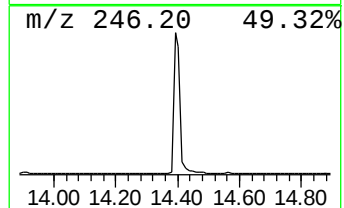
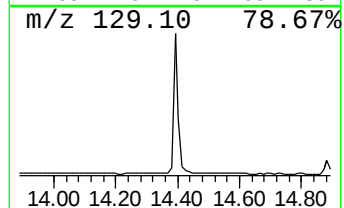
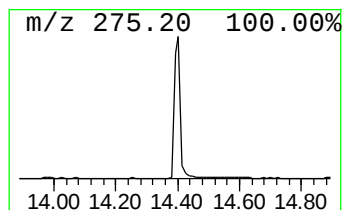
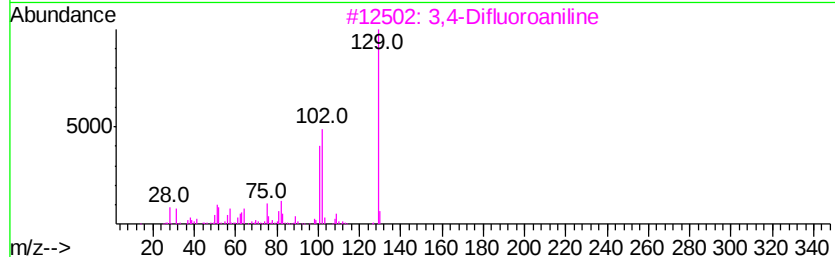
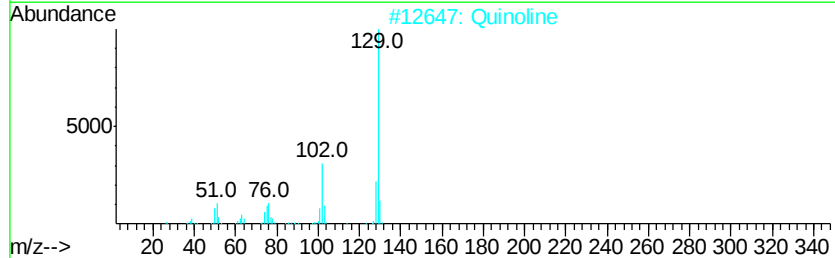
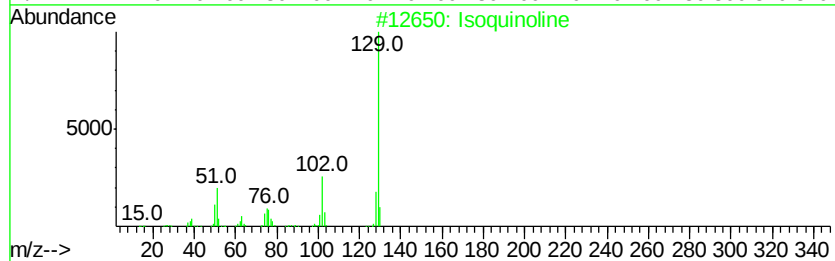
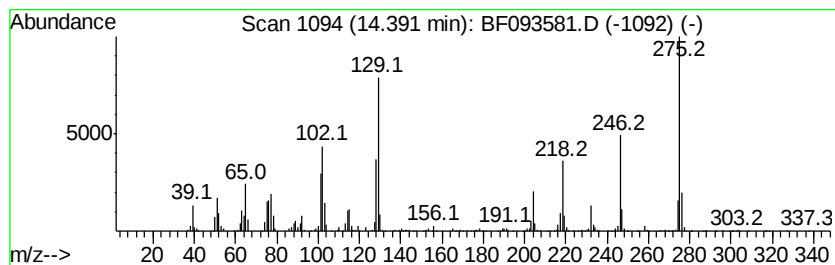
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 unknown14.39 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.39	19.39 ng	1185440	Chrysene-d12	13.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Isoquinoline	129	C9H7N	000119-65-3	38
2		Quinoline	129	C9H7N	000091-22-5	38
3		3,4-Difluoroaniline	129	C6H5F2N	003863-11-4	30
4		1-Methyl-5'-phenylspiro(indoline...	275	C17H13N3O	015970-21-5	30
5		Ethyl 2-cyano-2-(9-fluorenyliden...	275	C18H13N02	014003-25-9	25



Data Path : Z:\HPCHEM1\BNA F\DATA\BF031117\
Data File : BF093581.D
Acq On : 11 Mar 2017 21:38
Operator : SJ/MA
Sample : I2228-02 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TEC-TP3

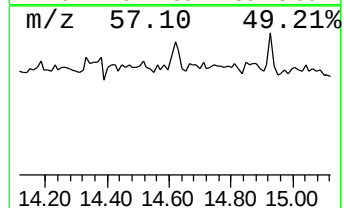
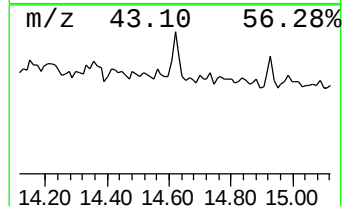
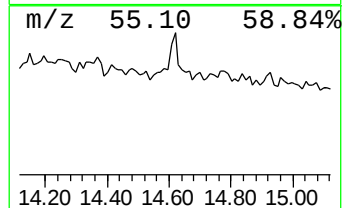
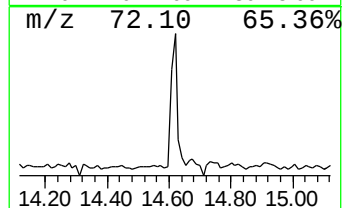
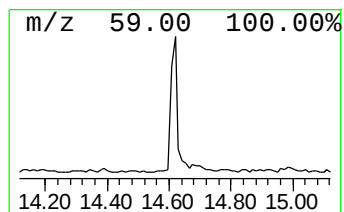
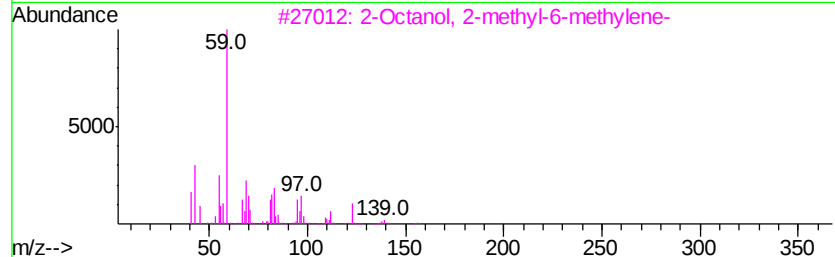
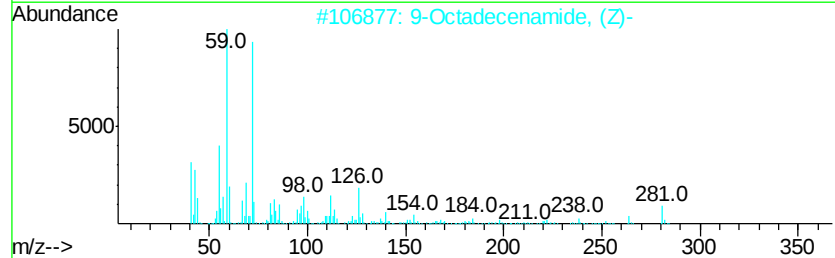
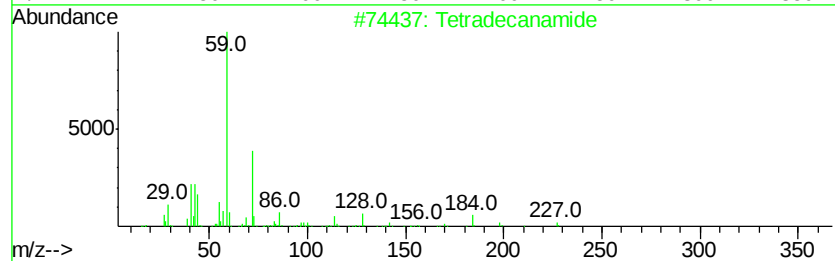
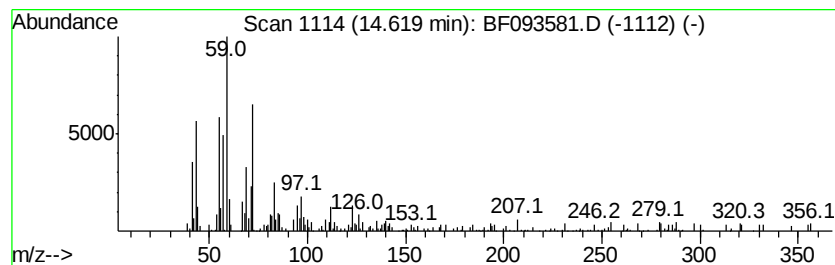
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Tetradecanamide Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.62	2.17 ng	139735	Perylene-d12	15.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetradecanamide	227	C14H29NO	000638-58-4	59
2		9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	59
3		2-Octanol, 2-methyl-6-methylene-	156	C10H20O	018479-59-9	50
4		Octadecanamide	283	C18H37NO	000124-26-5	38
5		Acetic acid, cyanohydroxyimino-,...	128	C4H4N2O3	061295-92-9	38



Data Path : Z:\HPCHEM1\BNA F\DATA\BF031117\
Data File : BF093581.D
Acq On : 11 Mar 2017 21:38
Operator : SJ/MA
Sample : I2228-02 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TEC-TP3

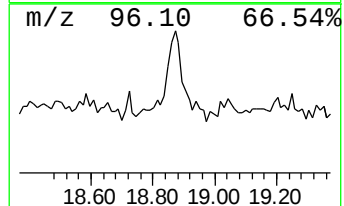
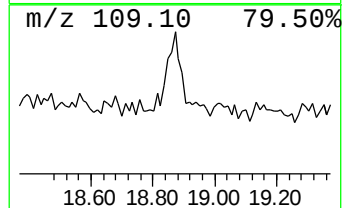
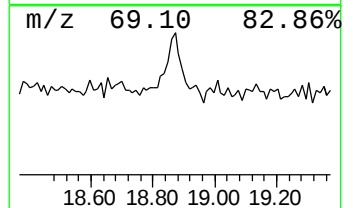
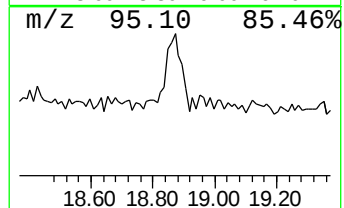
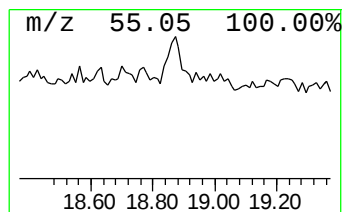
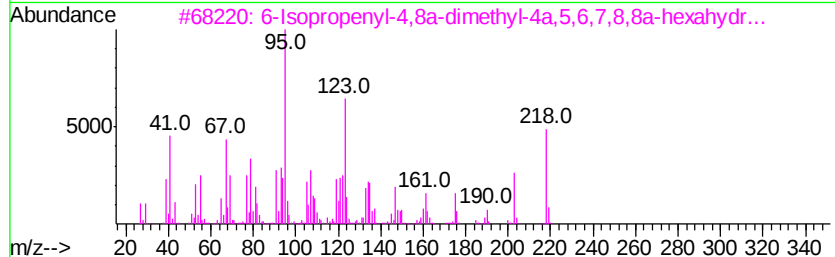
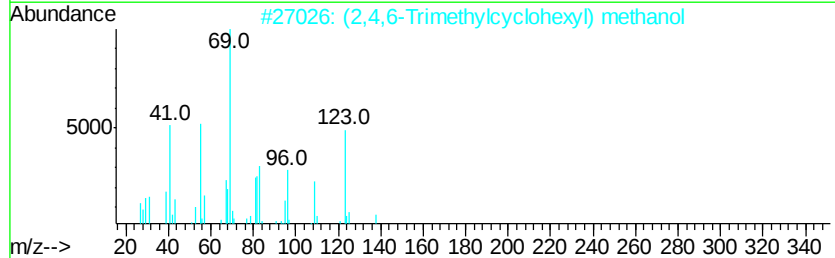
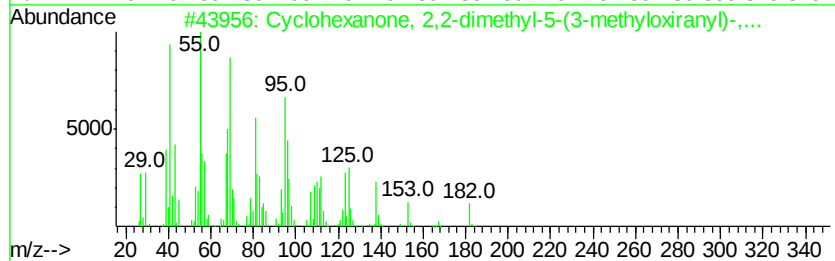
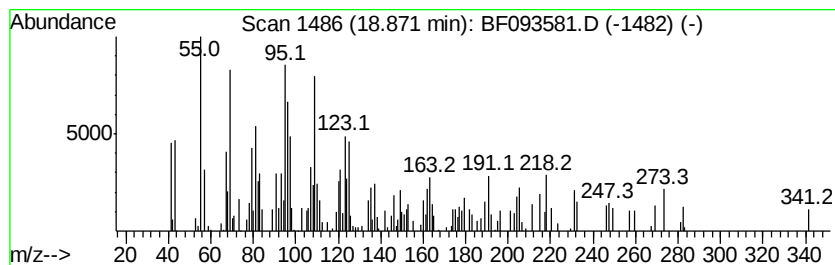
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF030717.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 unknown18.87 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.87	2.88 ng	185861	Perylene-d12	15.27

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexanone, 2,2-dimethyl-5-(3...	182	C11H18O2	141033-65-0	35
2			(2,4,6-Trimethylcyclohexyl) meth...	156	C10H20O	013702-56-2	22
3			6-Isopropenyl-4,8a-dimethyl-4a,5...	218	C15H22O	086917-79-5	18
4			5-.beta..beta.-dimethyl-4-methy...	196	C10H16N2O2	1000129-15-7	15
5			2,6,10-Nonadecatriene, 2,6,10,14...	318	C23H42	031147-35-0	14



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 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TEC-TP3

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF030717.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
2-Pentanone, 4-hy...	4.86	5.9	ng	493015	1	6.72	1677400	20.0
unknown6.49	6.49	30.0	ng	2515050	1	6.72	1677400	20.0
Octane, 2,6-dimet...	8.42	3.4	ng	357805	2	8.00	2105150	20.0
Octane, 2-methyl-	9.02	3.3	ng	311868	3	9.75	1886460	20.0
Decahydro-4,4,8,9...	9.14	4.8	ng	451768	3	9.75	1886460	20.0
unknown9.64	9.64	7.8	ng	740893	3	9.75	1886460	20.0
unknown10.01	10.01	2.4	ng	221725	3	9.75	1886460	20.0
2-Bromo dodecane	10.41	4.6	ng	433094	3	9.75	1886460	20.0
Dodecane, 2-methy...	10.68	10.1	ng	832340	4	11.24	1652160	20.0
Tridecane	11.13	5.2	ng	432623	4	11.24	1652160	20.0
unknown14.39	14.39	19.4	ng	1185440	5	13.88	1222980	20.0
Tetradecanamide	14.62	2.2	ng	139735	6	15.27	1288560	20.0
unknown18.87	18.87	2.9	ng	185861	6	15.27	1288560	20.0