

Data Path : Z:\HPCHEM1\BNA F\DATA\BF032718\
 Data File : BF103979.D
 Acq On : 27 Mar 2018 19:29
 Operator : JU/SJ
 Sample : J2065-01
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RO-CV-2536

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-BF031518.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	2.316	33	39	46	rVB	103886	142174	4.70%	0.403%
2	5.222	528	533	545	rBV	72395	117878	3.90%	0.334%
3	5.610	594	599	628	rBV	1983734	2884270	95.34%	8.174%
4	6.604	763	768	788	rBV	2175026	2998078	99.11%	8.497%
5	6.745	788	792	808	rVB	2729977	2983816	98.63%	8.456%
6	6.969	827	830	846	rBV	713990	746554	24.68%	2.116%
7	7.128	853	857	878	rVB	2302200	2217294	73.30%	6.284%
8	7.534	922	926	935	rBV	1523694	1750611	57.87%	4.961%
9	8.251	1045	1048	1054	rBV	939177	950934	31.43%	2.695%
10	9.204	1204	1210	1214	rBV7	70263	171420	5.67%	0.486%
11	9.239	1214	1216	1221	rVV3	120151	149063	4.93%	0.422%
12	9.328	1225	1231	1235	rBV	3144927	3025118	100.00%	8.573%
13	9.398	1239	1243	1246	rBV2	359332	377734	12.49%	1.071%
14	9.633	1281	1283	1285	rBV2	130393	144205	4.77%	0.409%
15	9.657	1285	1287	1292	rVV3	236361	374641	12.38%	1.062%
16	9.716	1293	1297	1300	rBV2	1938047	1716634	56.75%	4.865%
17	9.745	1300	1302	1304	rVB2	192754	152695	5.05%	0.433%
18	9.769	1304	1306	1310	rBV2	496381	562513	18.59%	1.594%
19	9.804	1310	1312	1313	rBV2	148087	103869	3.43%	0.294%
20	9.875	1317	1324	1326	rBV4	1345636	1589381	52.54%	4.504%
21	9.898	1326	1328	1329	rBV2	280530	250720	8.29%	0.711%
22	9.922	1329	1332	1335	rVB	2089265	1855198	61.33%	5.258%
23	9.963	1335	1339	1340	rBV2	605274	657373	21.73%	1.863%
24	9.992	1341	1344	1347	rBV3	916142	1350875	44.66%	3.828%
25	10.045	1350	1353	1354	rBV3	444585	453065	14.98%	1.284%
26	10.151	1369	1371	1376	rBV2	429101	771394	25.50%	2.186%
27	10.257	1387	1389	1393	rBV3	456172	643808	21.28%	1.825%
28	10.333	1399	1402	1403	rBV2	880430	857222	28.34%	2.429%
29	10.428	1414	1418	1421	rBV4	1832130	2087928	69.02%	5.917%
30	13.104	1869	1873	1879	rVB	1611876	1730763	57.21%	4.905%
31	14.163	2049	2053	2056	rVB	738256	660945	21.85%	1.873%
32	14.915	2178	2181	2184	rBV	129783	128283	4.24%	0.364%
33	15.074	2206	2208	2212	rVB4	58479	59742	1.97%	0.169%
34	15.698	2309	2314	2322	rVB	473851	618801	20.46%	1.754%

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Sample : J2065-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
RO-CV-2536

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-BF031518.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

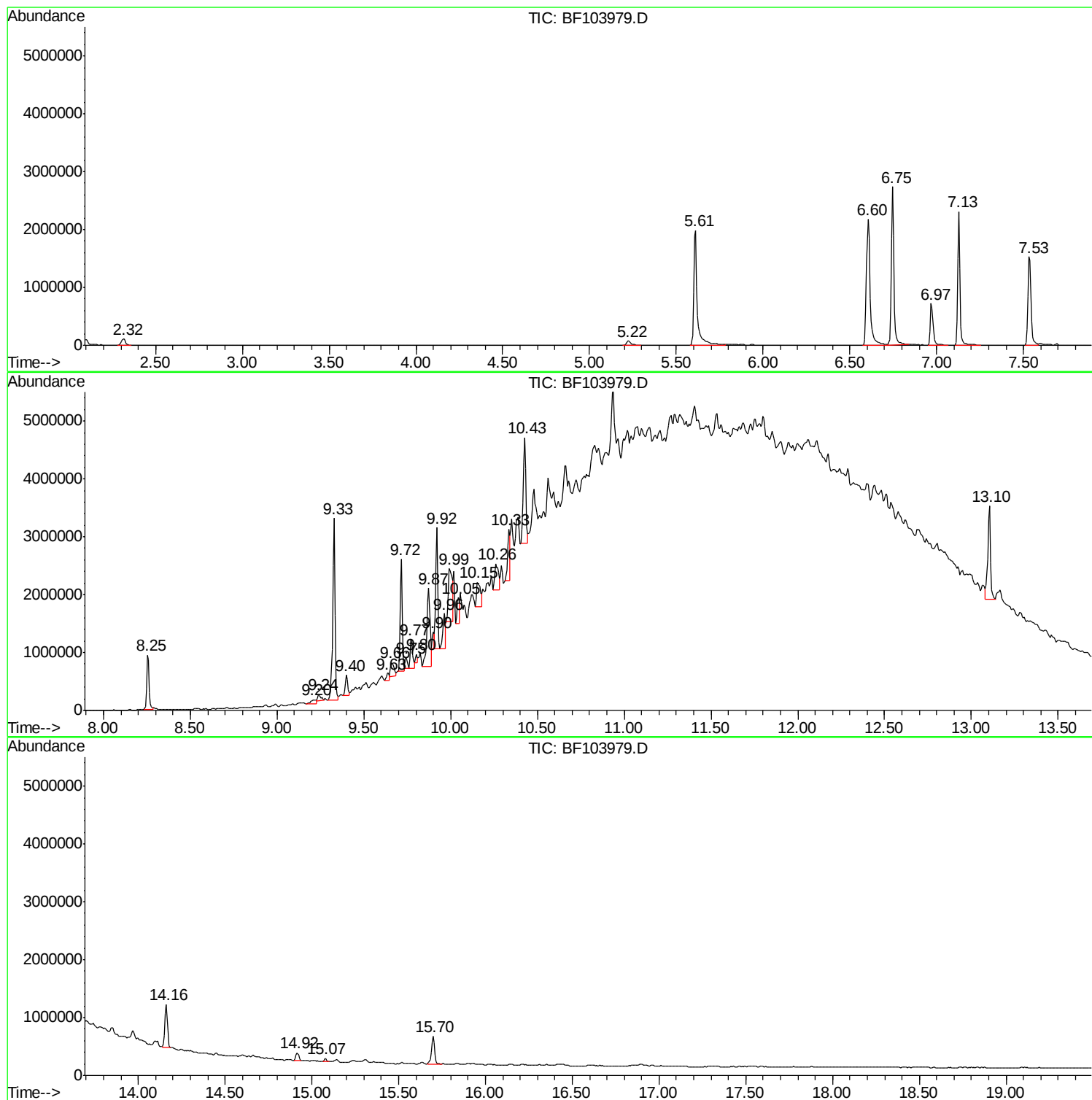
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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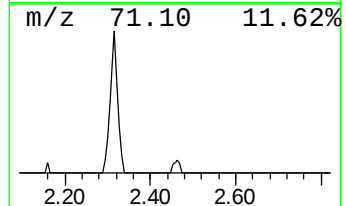
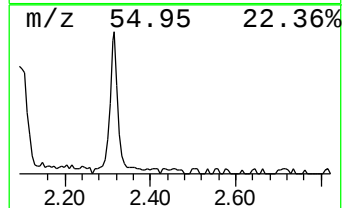
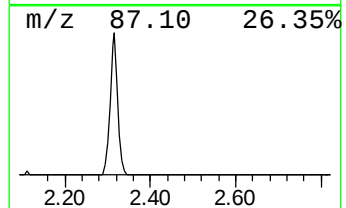
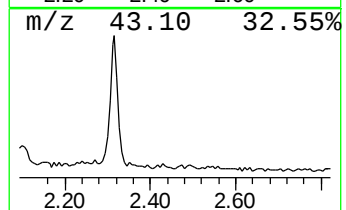
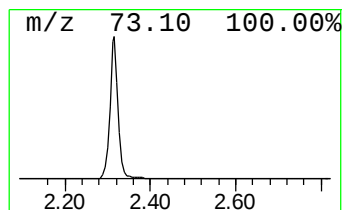
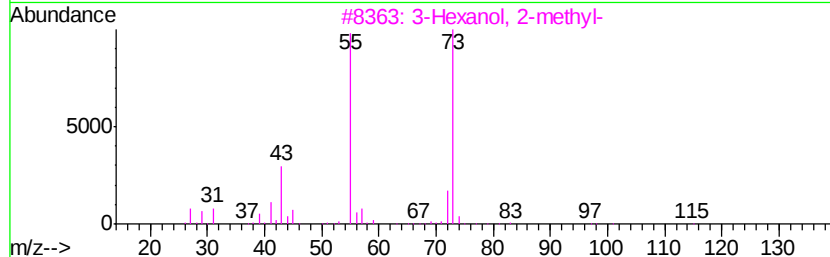
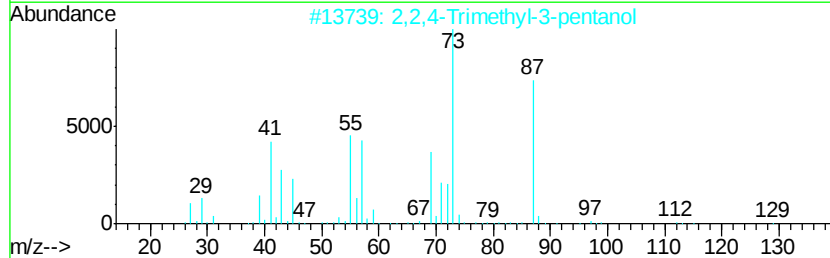
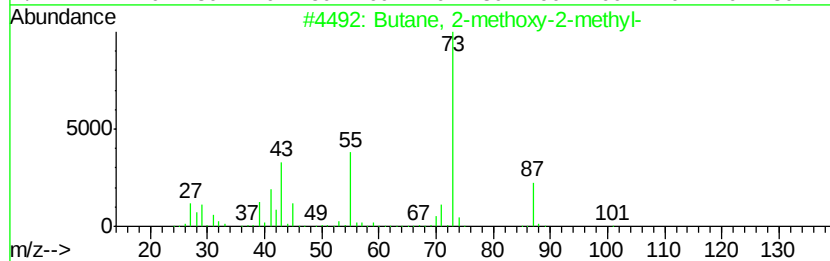
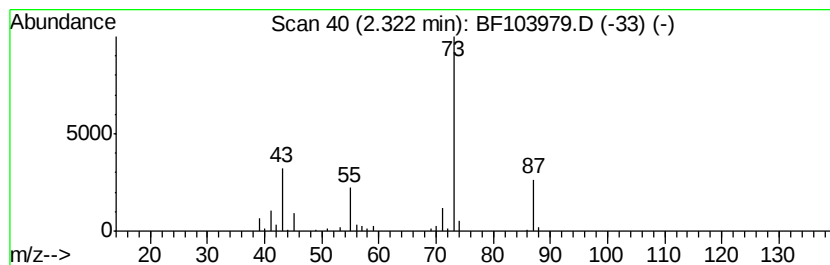
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.32	3.81 ng	142174	1,4-Dichlorobenzene-d4	6.97

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
3			3-Hexanol, 2-methyl-	116	C7H16O	000617-29-8	17
4			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17
5			N-Ethylformamide	73	C3H7NO	000627-45-2	9



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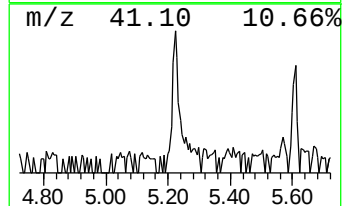
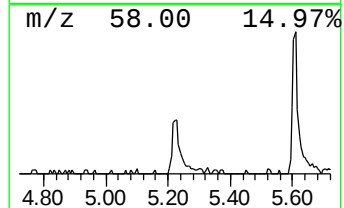
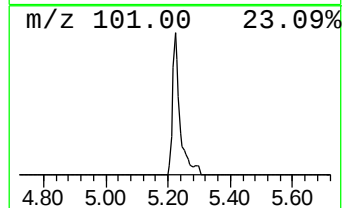
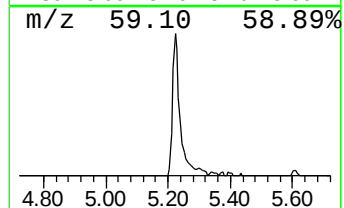
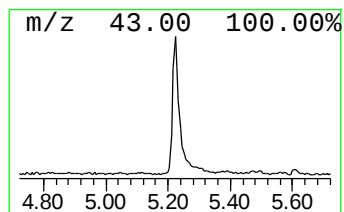
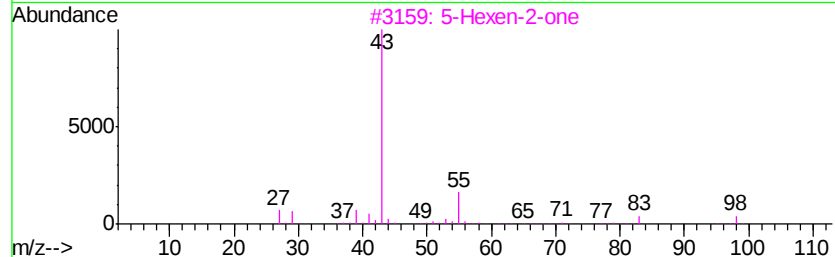
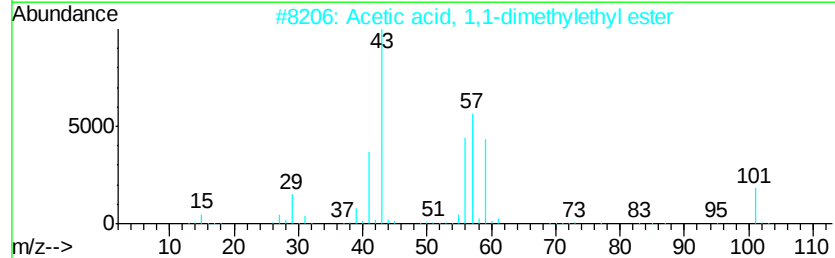
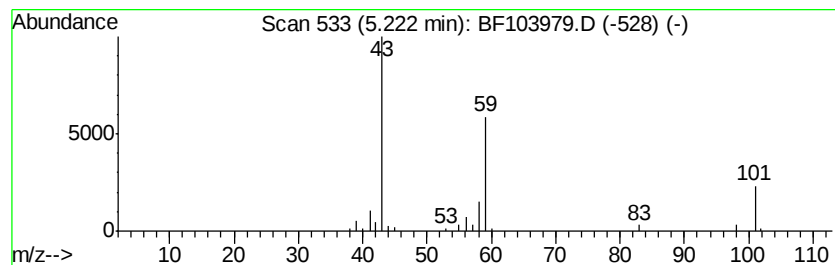
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.22	3.16 ng	117878	1,4-Dichlorobenzene-d4	6.97

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	38
3			5-Hexen-2-one	98	C6H10O	000109-49-9	9
4			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5			2-Butanone, 3-ethoxy-3-methyl-	130	C7H14O2	036687-99-7	9



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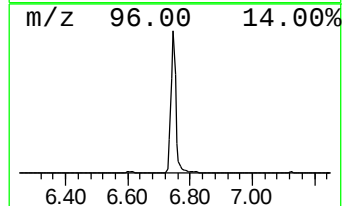
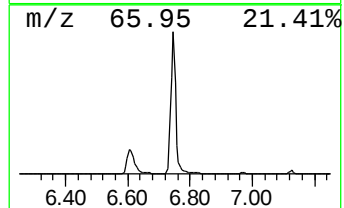
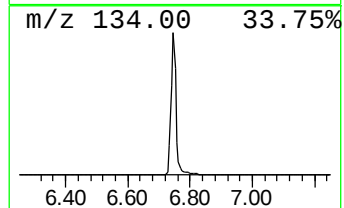
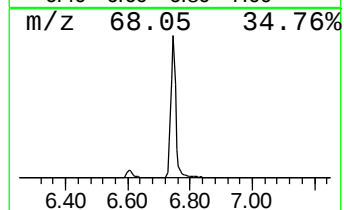
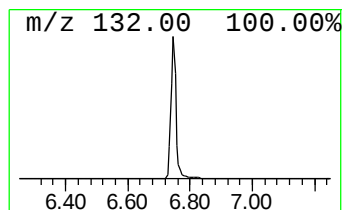
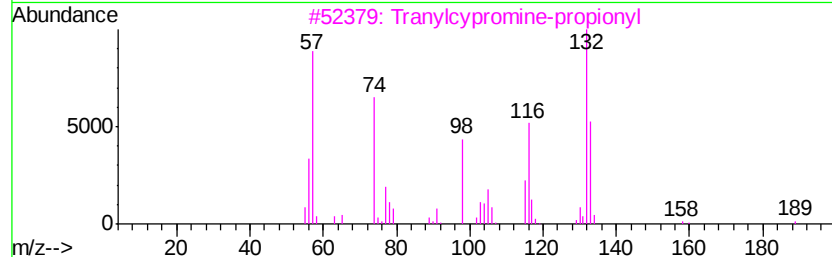
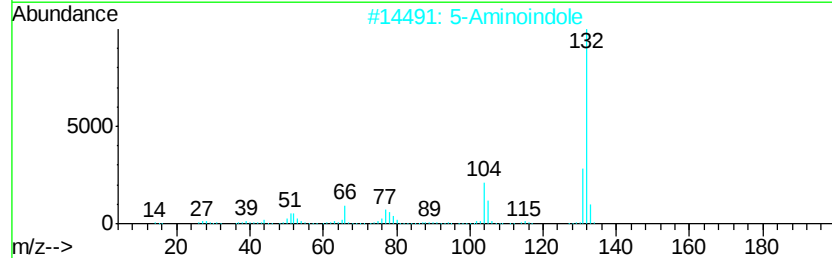
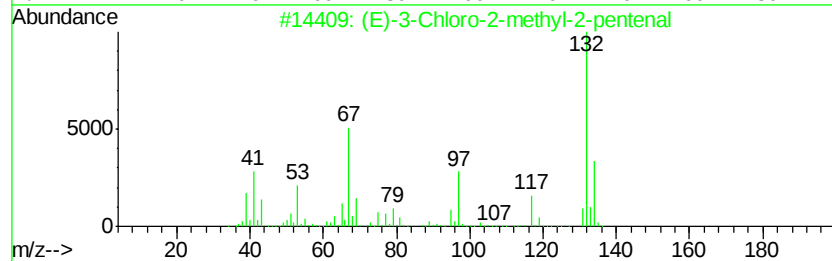
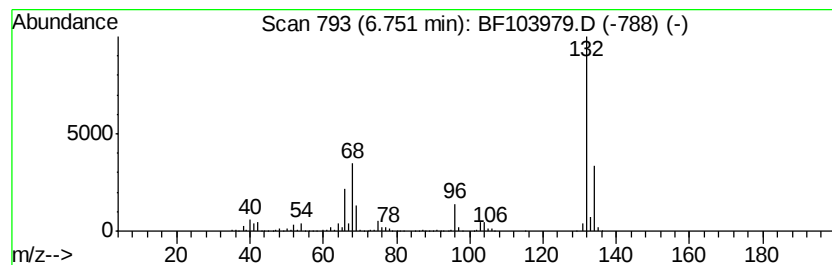
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown6.75 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.75	79.94 ng	2983820	1,4-Dichlorobenzene-d4	6.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
2		5-Aminoindole	132	C8H8N2	005192-03-0	12
3		Tranvlcyproline-propionyl	189	C12H15NO	1000123-86-3	10
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9
5		4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9



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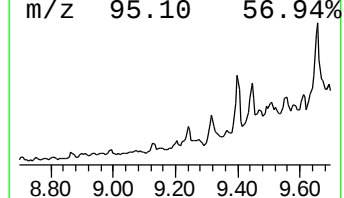
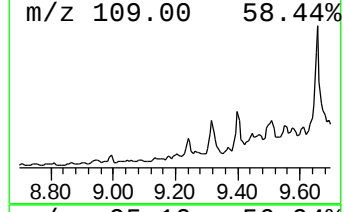
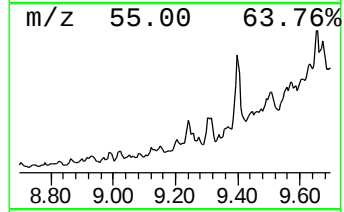
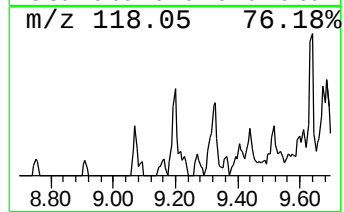
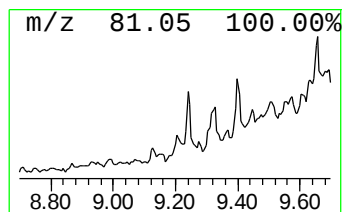
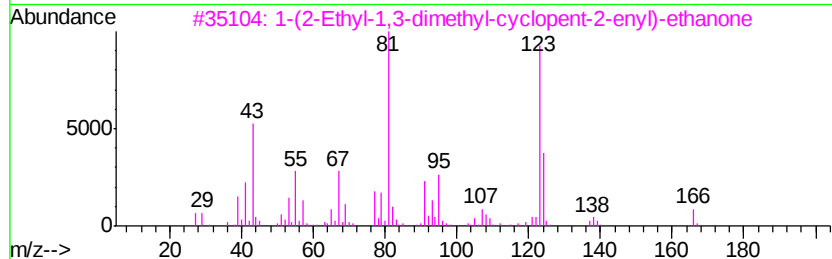
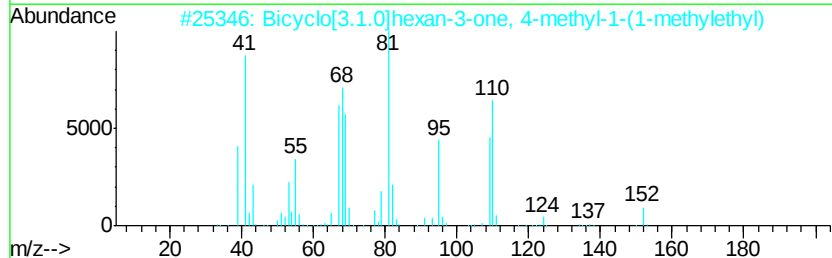
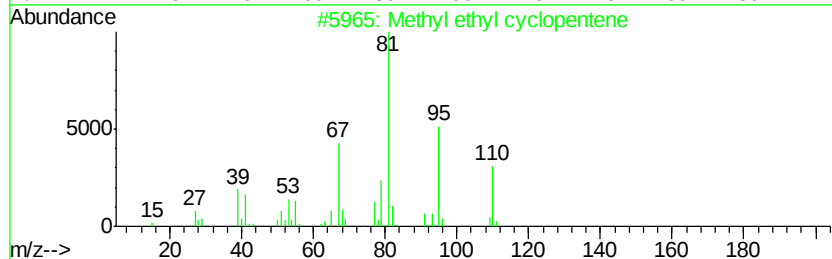
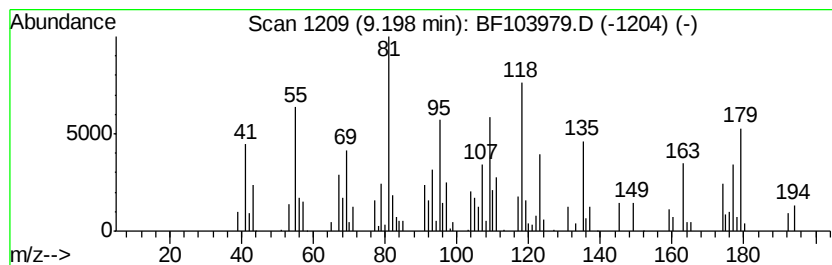
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 unknown9.20 Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.20	2.54 ng	171420	Acenaphthene-d10	10.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Methyl ethyl cyclopentene	110	C8H14	019780-56-4	38
2		Bicyclo[3.1.0]hexan-3-one, 4-met...	152	C10H16O	001125-12-8	35
3		1-(2-Ethyl-1,3-dimethyl-cyclopent...	166	C11H18O	1000185-18-1	27
4		trans-(2-Ethylcyclopentyl)methanol	128	C8H16O	036258-08-9	27
5		2,5,9-Tetradecatriene, 3,12-diet...	248	C18H32	074685-87-3	27



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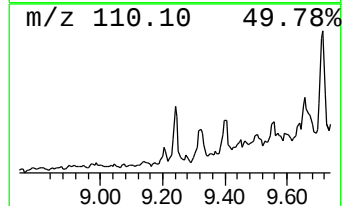
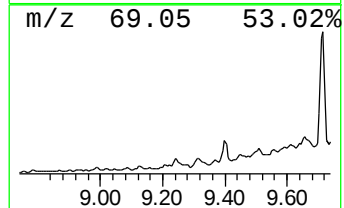
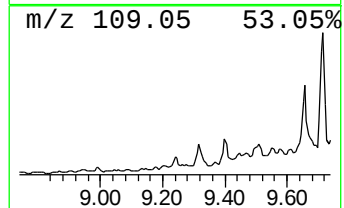
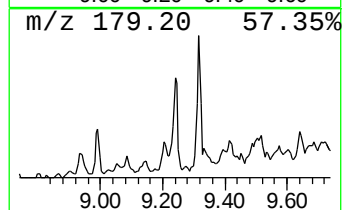
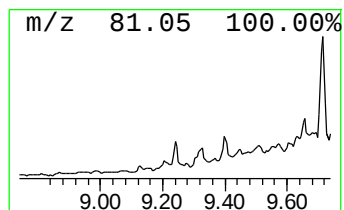
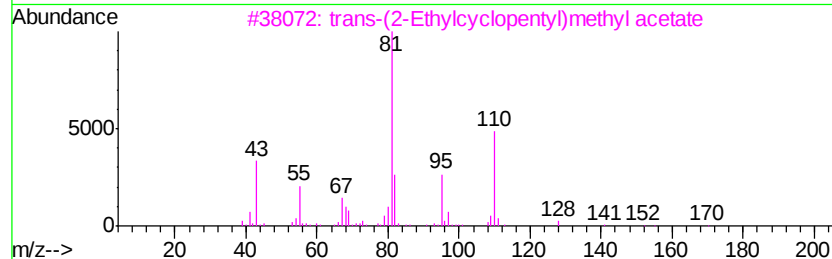
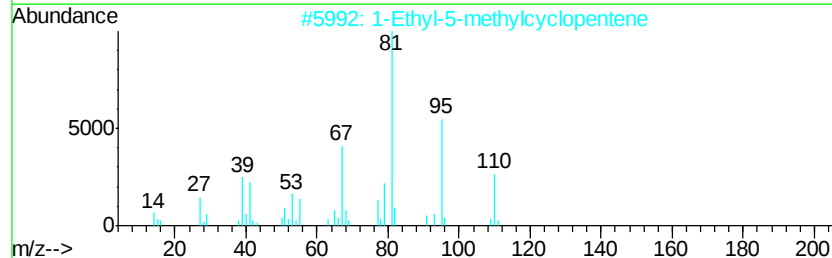
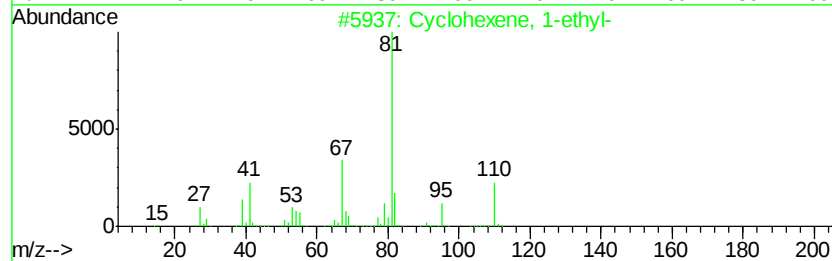
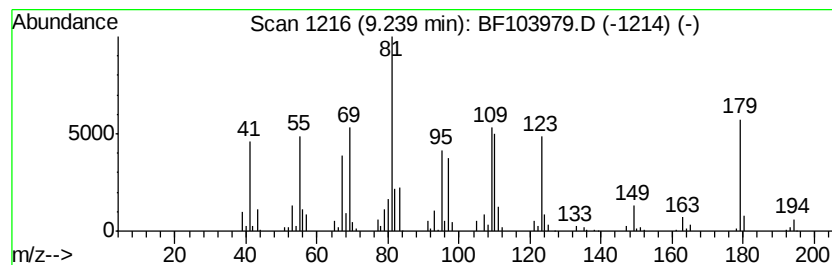
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 unknown9.24 Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.24	2.21 ng	149063	Acenaphthene-d10	10.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexene, 1-ethyl-	110	C8H14	001453-24-3	38
2		1-Ethyl-5-methylcyclopentene	110	C8H14	097797-57-4	35
3		trans-(2-Ethylcyclopentyl)methyl...	170	C10H18O2	1000099-13-9	35
4		4-Ethylcyclohexene	110	C8H14	003742-42-5	27
5		Cyclohexene, 3-ethyl-	110	C8H14	002808-71-1	27



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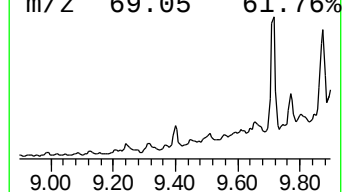
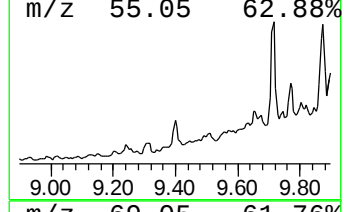
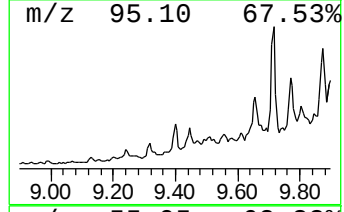
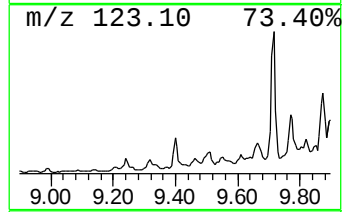
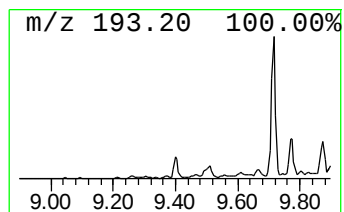
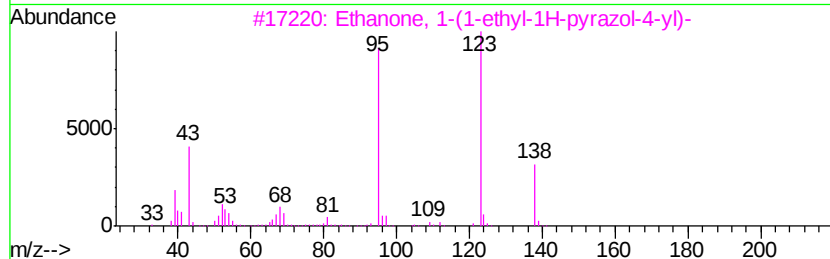
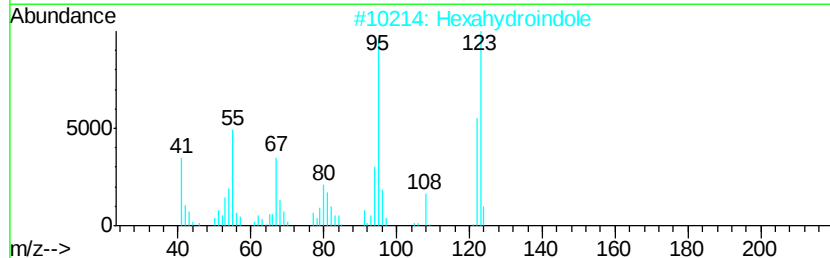
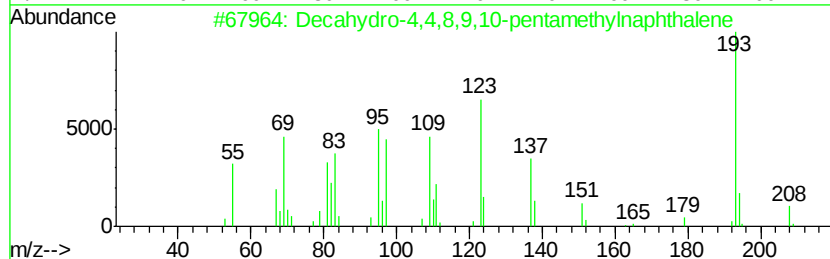
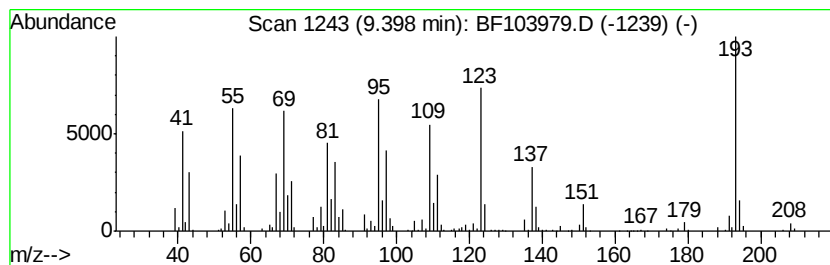
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Decahydro-4,4,8,9,10-pentam... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.40	5.59 ng	377734	Acenaphthene-d10	10.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Decahydro-4,4,8,9,10-pentamethyl...	208	C15H28	080655-44-3	91
2		Hexahydroindole	123	C8H13N	018159-32-5	30
3		Ethanone, 1-(1-ethyl-1H-pyrazol-...	138	C7H10N2O	1000316-95-0	25
4		Cyclopentene, 1,3-dimethyl-2-(1-...	138	C10H18	061142-32-3	25
5		1,3-Benzenediamine, 4-methoxy-	138	C7H10N2O	000615-05-4	22



Data Path : Z:\HPCHEM1\BNA F\DATA\BF032718\
Data File : BF103979.D
Acq On : 27 Mar 2018 19:29
Operator : JU/SJ
Sample : J2065-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
RO-CV-2536

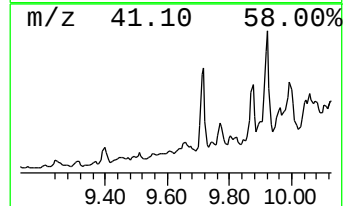
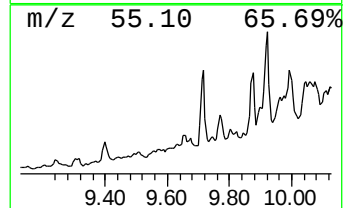
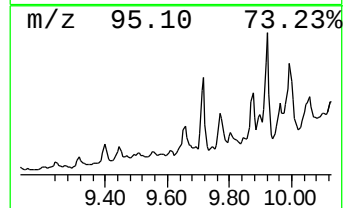
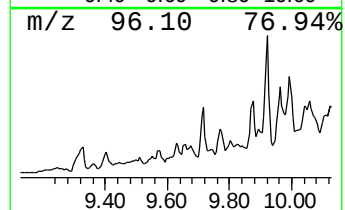
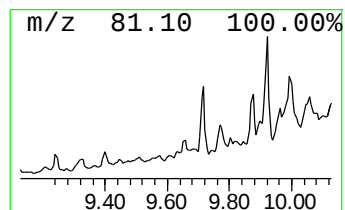
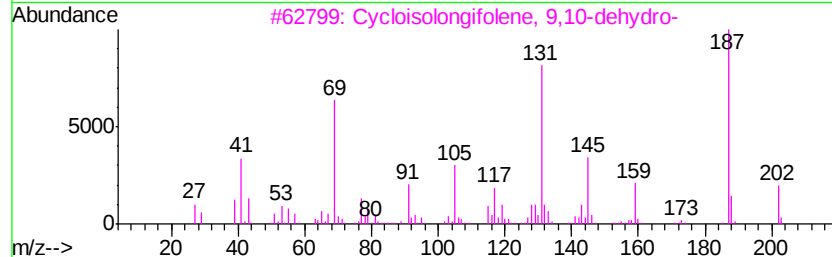
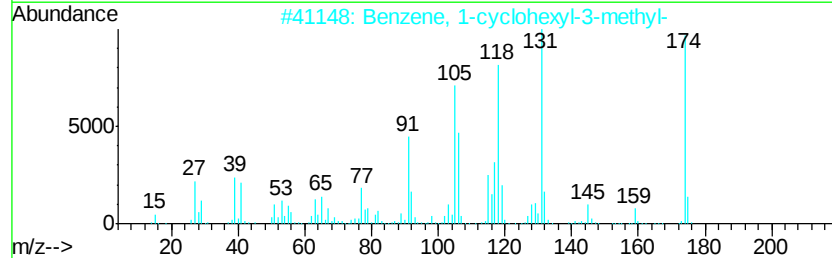
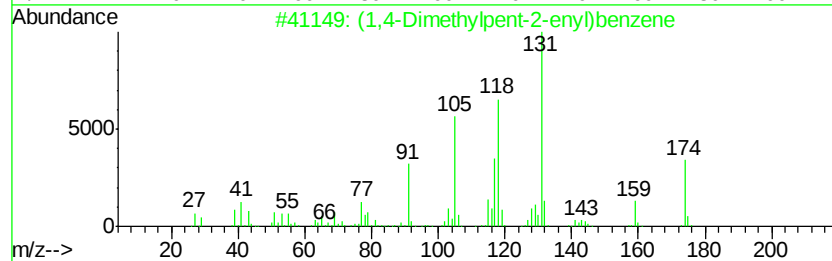
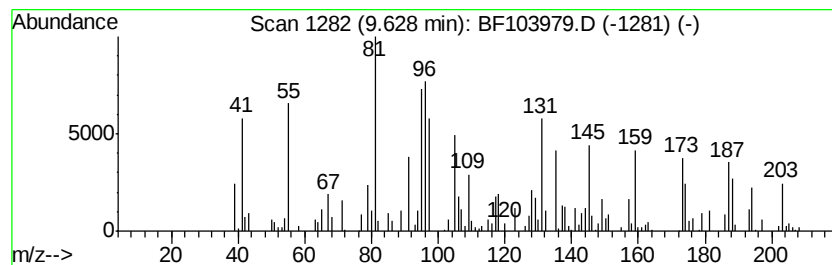
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 unknown9.63 Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.63	2.13 ng	144205	Acenaphthene-d10	10.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(1,4-Dimethylpent-2-enyl)benzene	174	C13H18	1000184-97-9	46
2		Benzene, 1-cyclohexyl-3-methyl-	174	C13H18	004575-46-6	41
3		Cycloisolongifolene, 9,10-dehydro-	202	C15H22	1000156-81-6	30
4		1H-1,3,2-Benzodiazaborole, 2-but...	174	C10H15BN2	031748-14-8	25
5		3-Fluorobenzoic acid, 3-phenylpr...	258	C16H15FO2	1000279-05-0	22



Data Path : Z:\HPCHEM1\BNA F\DATA\BF032718\
Data File : BF103979.D
Acq On : 27 Mar 2018 19:29
Operator : JU/SJ
Sample : J2065-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
RO-CV-2536

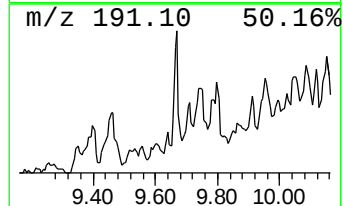
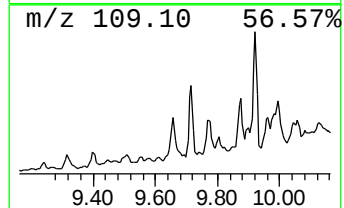
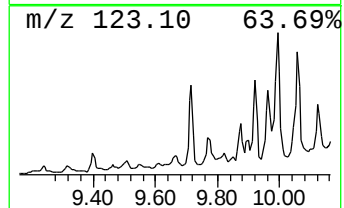
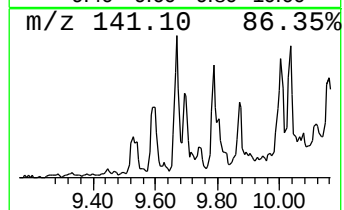
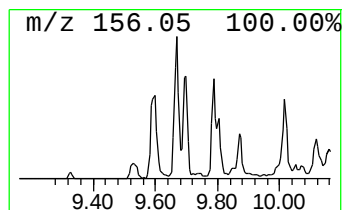
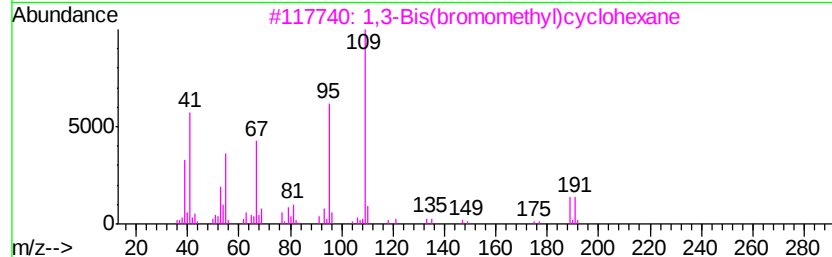
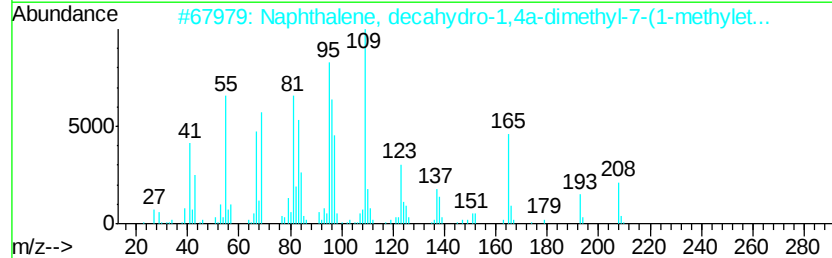
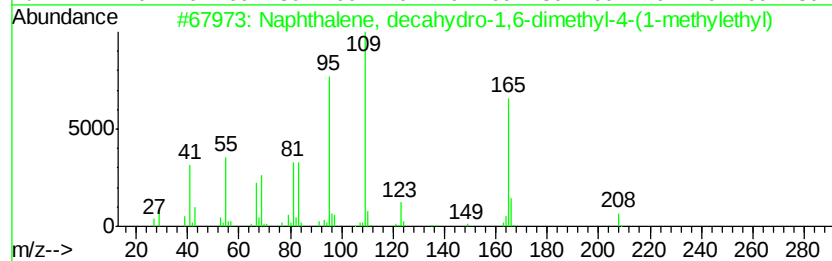
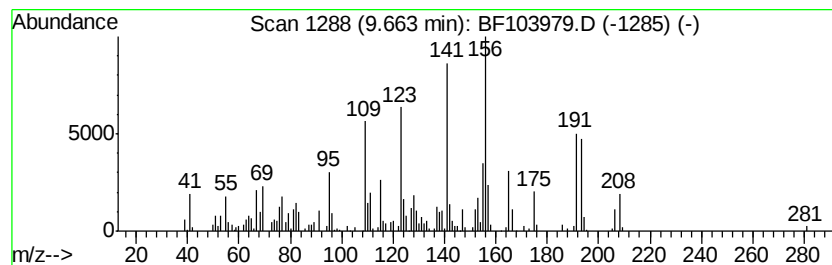
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Naphthalene, decahydro-1,6-... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.66	5.55 ng	374641	Acenaphthene-d10	10.02

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, decahydro-1,6-dimet...	208	C15H28	029788-41-8	74
2			Naphthalene, decahydro-1,4a-dime...	208	C15H28	030824-81-8	43
3			1,3-Bis(bromomethyl)cyclohexane	268	C8H14Br2	1000216-89-9	43
4			1-Methoxyadamantane	166	C11H18O	006221-74-5	35
5			Tricyclo[4.3.1.1(3,8)]undecan-3-ol	166	C11H18O	014504-80-4	30



Data Path : Z:\HPCHEM1\BNA F\DATA\BF032718\
Data File : BF103979.D
Acq On : 27 Mar 2018 19:29
Operator : JU/SJ
Sample : J2065-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampled :
RO-CV-2536

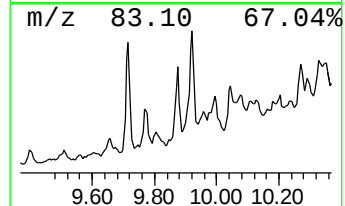
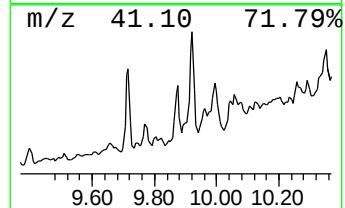
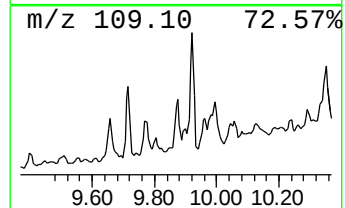
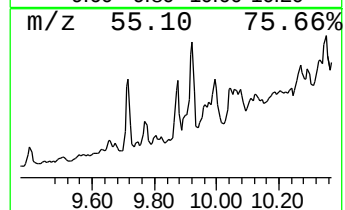
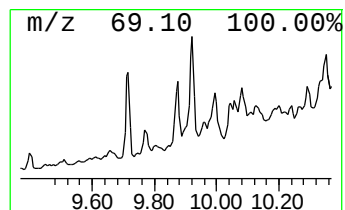
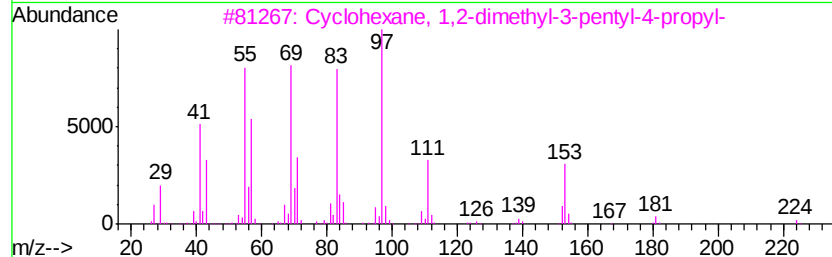
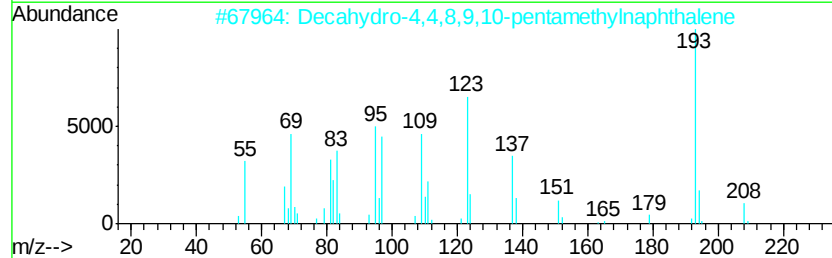
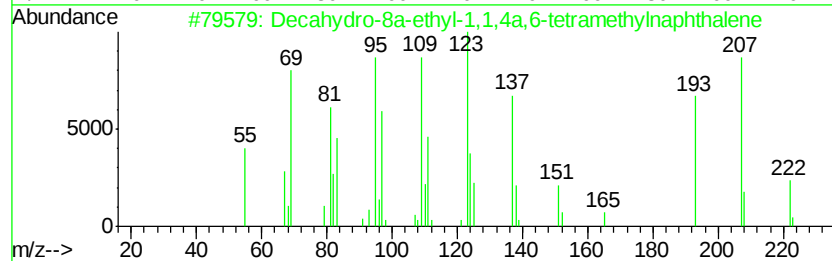
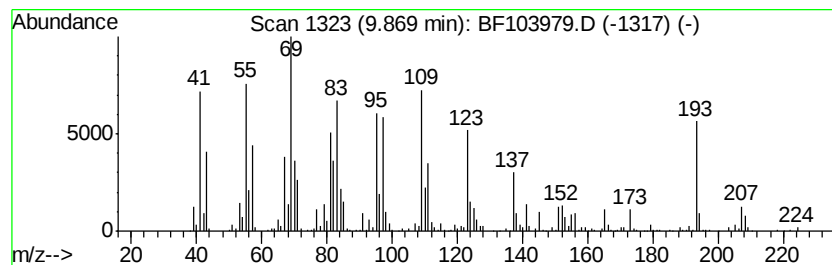
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Decahydro-8a-ethyl-1,1,4a,6... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.87	23.53 ng	1589380	Acenaphthene-d10	10.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Decahydro-8a-ethyl-1,1,4a,6-tetr...	222	C16H30	1000100-23-6	72
2		Decahydro-4,4,8,9,10-pentamethyl...	208	C15H28	080655-44-3	64
3		Cyclohexane, 1,2-dimethyl-3-pent...	224	C16H32	062376-17-4	41
4		Bicyclo[3.1.1]heptan-3-one, 2,6,...	152	C10H16O	015358-88-0	38
5		Bicyclo[3.1.1]heptan-3-one, 2,6,...	152	C10H16O	000547-60-4	38



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032718\
Data File : BF103979.D
Acq On : 27 Mar 2018 19:29
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Misc :
ALS Vial : 11 Sample Multiplier: 1

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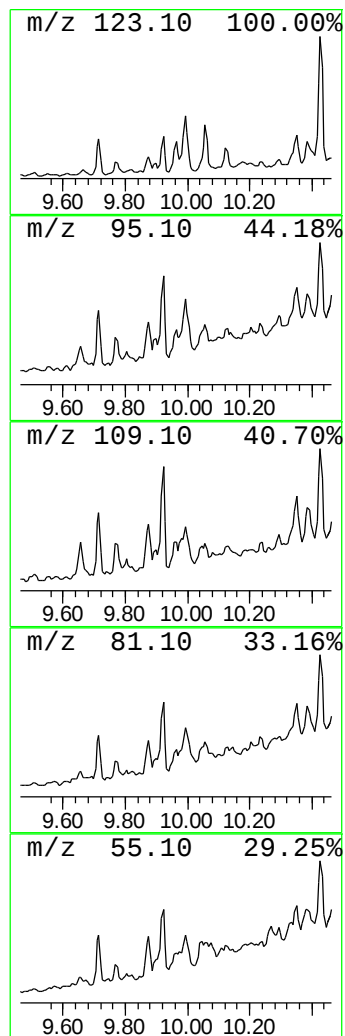
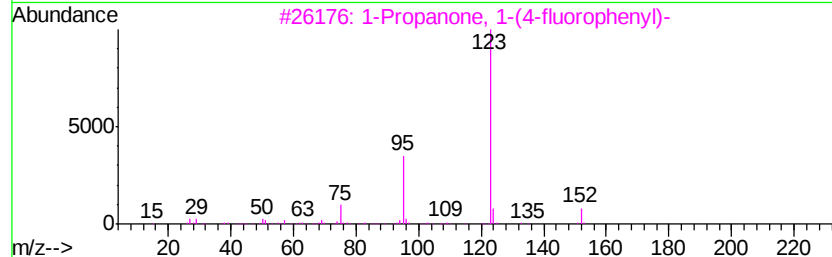
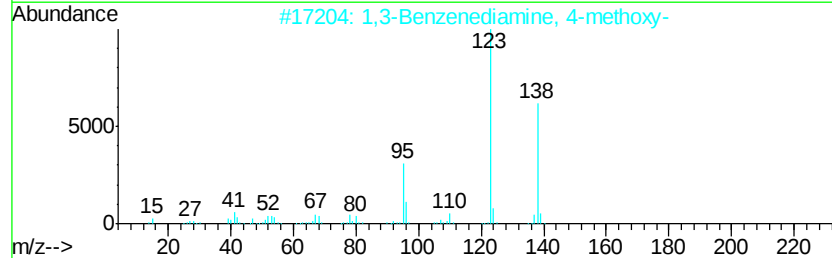
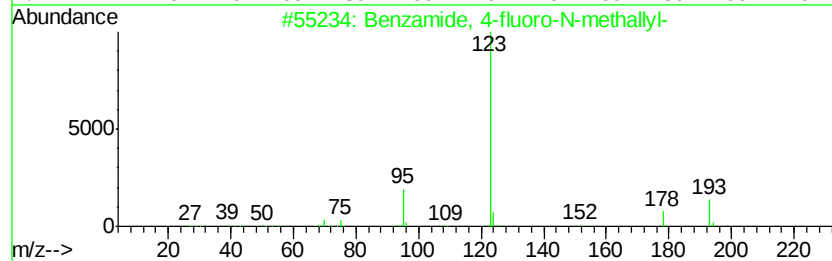
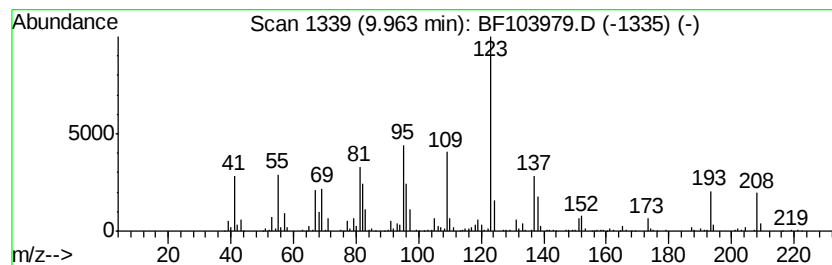
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 unknown9.96 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.96	9.73 ng	657373	Acenaphthene-d10	10.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzamide, 4-fluoro-N-methallyl-	193	C11H12FNO	1000340-31-9	43
2		1,3-Benzenediamine, 4-methoxy-	138	C7H10N2O	000615-05-4	43
3		1-Propanone, 1-(4-fluorophenyl)-	152	C9H9FO	000456-03-1	43
4		Imidazolidin-2-one, 1-(2-fluorob...	208	C10H9FN2O2	1000310-62-2	43
5		Ethanone, 1-(4-fluorophenyl)-	138	C8H7FO	000403-42-9	43



Data Path : Z:\HPCHEM1\BNA F\DATA\BF032718\
Data File : BF103979.D
Acq On : 27 Mar 2018 19:29
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RO-CV-2536

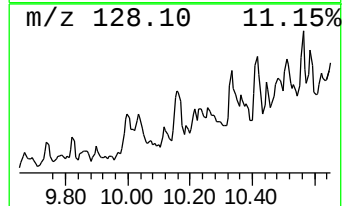
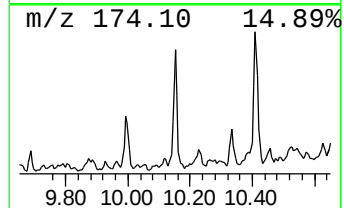
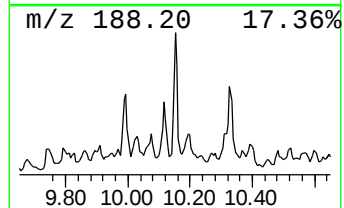
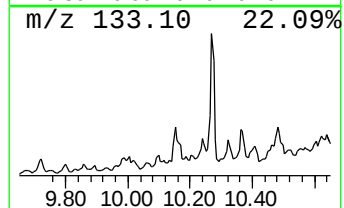
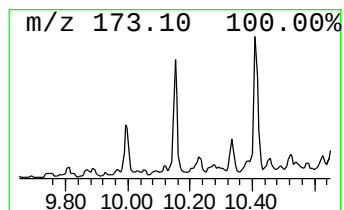
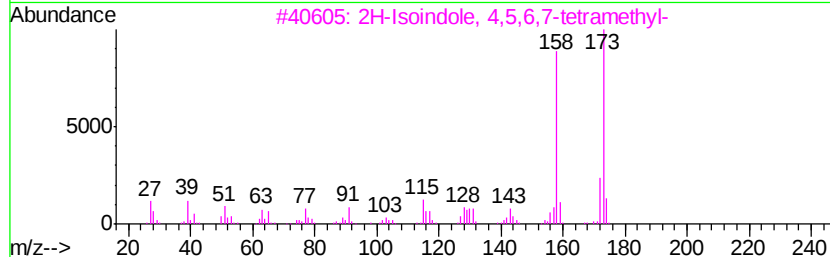
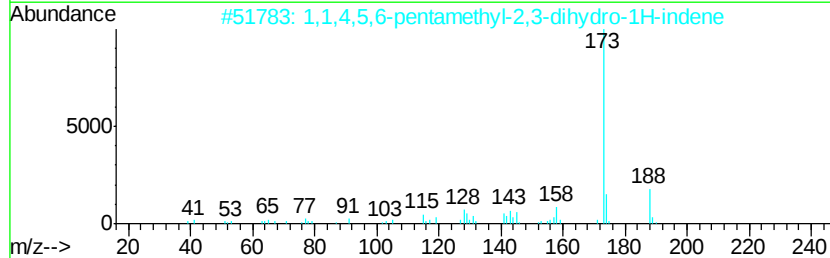
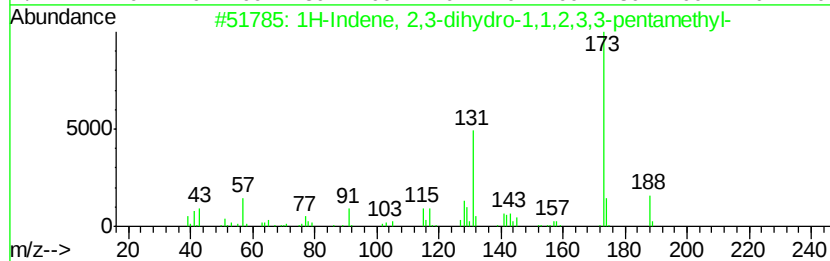
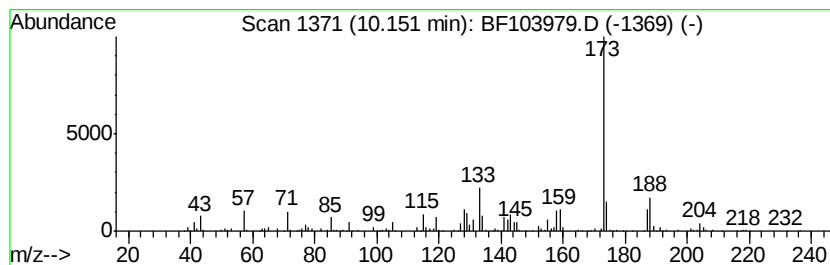
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 1H-Indene, 2,3-dihydro-1,1,... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.15	11.42 ng	771394	Acenaphthene-d10	10.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2,3-dihydro-1,1,2,3,3...	188	C14H20	001203-17-4	83
2		1,1,4,5,6-pentamethyl-2,3-dihydro...	188	C14H20	1000374-13-8	76
3		2H-Isoindole, 4,5,6,7-tetramethyl-	173	C12H15N	070187-61-0	64
4		4H-1-Benzopyran, 4,4,5,8-tetrame...	188	C13H16O	082391-06-8	50
5		1H-Inden-1-one, 2,3-dihydro-3,3,...	188	C13H16O	056298-98-7	50



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032718\
Data File : BF103979.D
Acq On : 27 Mar 2018 19:29
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ALS Vial : 11 Sample Multiplier: 1

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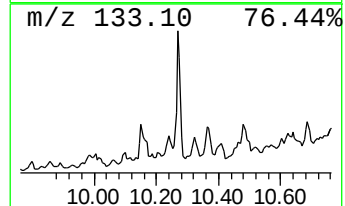
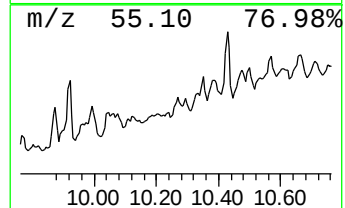
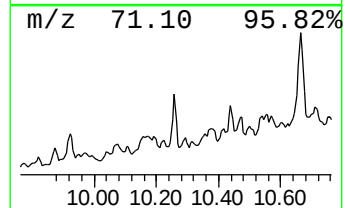
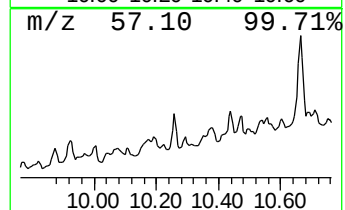
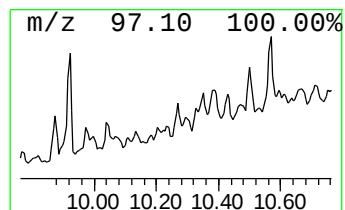
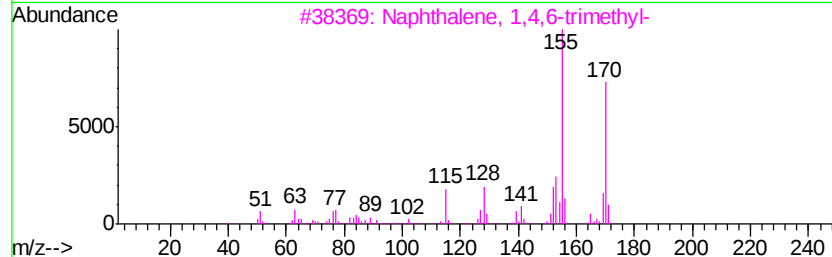
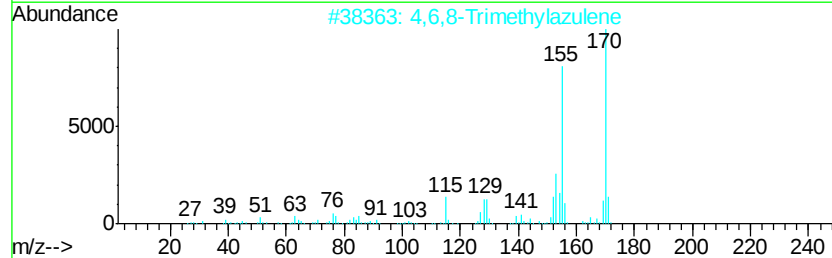
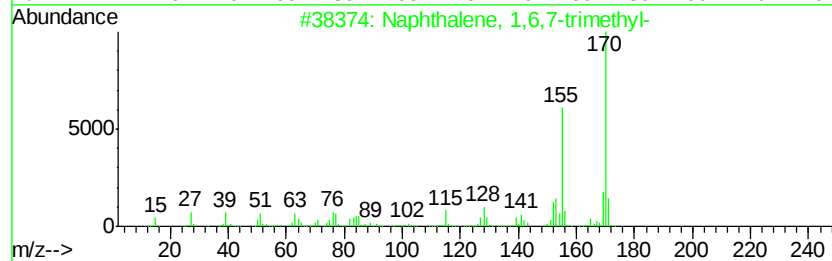
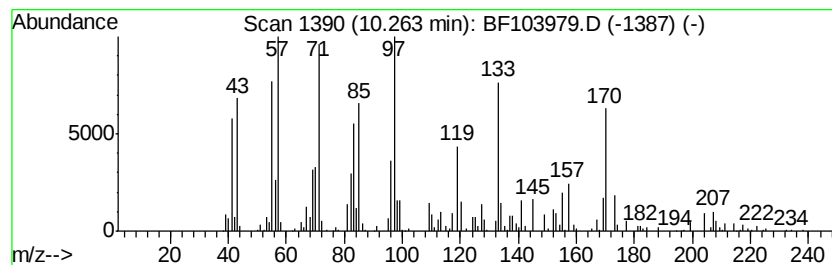
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Naphthalene, 1,6,7-trimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.26	9.53 ng	643808	Acenaphthene-d10	10.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	58
2		4,6,8-Trimethylazulene	170	C13H14	000941-81-1	43
3		Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	38
4		Octane, 2,4,6-trimethyl-	156	C11H24	062016-37-9	35
5		Nonane, 1-iodo-	254	C9H19I	004282-42-2	35



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032718\
Data File : BF103979.D
Acq On : 27 Mar 2018 19:29
Operator : JU/SJ
Sample : J2065-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

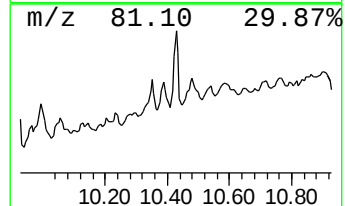
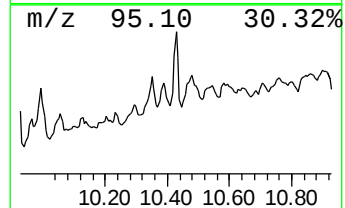
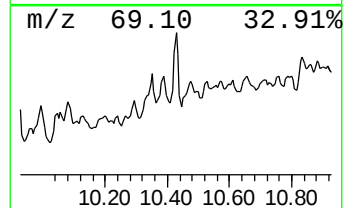
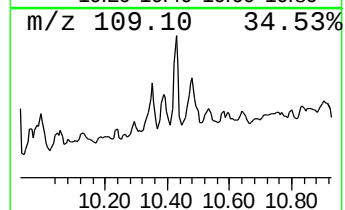
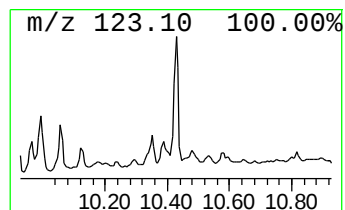
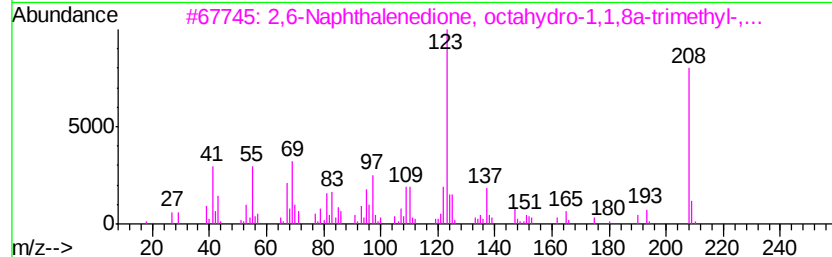
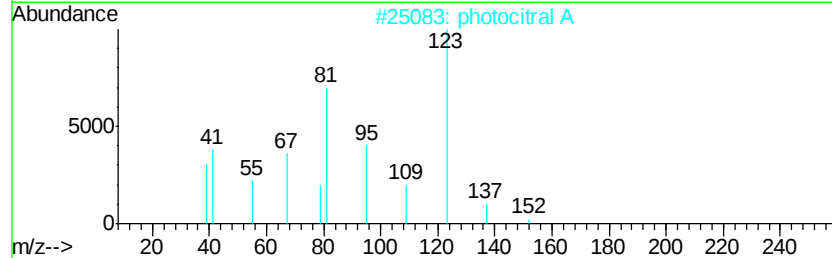
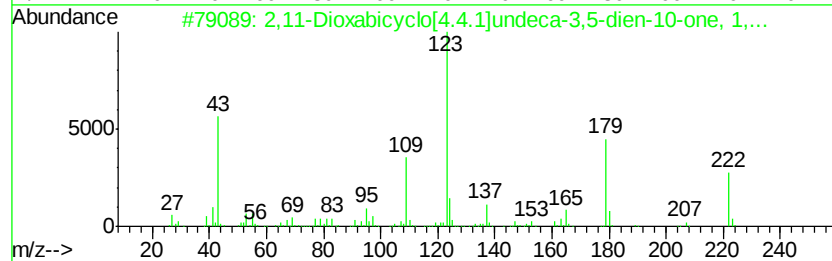
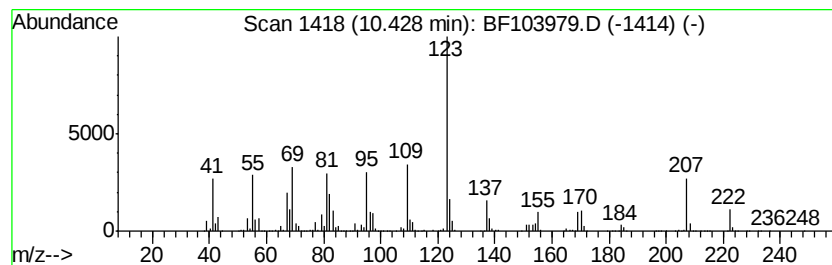
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ClientSampleId :
RO-CV-2536

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 2,11-Dioxabicyclo[4.4.1]und... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.43	30.91 ng	2087930	Acenaphthene-d10	10.02	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,11-Dioxabicyclo[4.4.1]undeca-3...	222	C13H18O3	070412-52-1	50
2	photocitral A	152	C10H16O	055253-28-6	47
3	2,6-Naphthalenedione, octahydro-...	208	C13H20O2	057289-16-4	43
4	3-Fluorobenzoic acid, 3-methylbu...	208	C12H13FO2	154559-38-3	43
5	3-Fluorobenzoic acid, 1-cyclopen...	236	C14H17FO2	1000279-04-7	38



Data Path : Z:\HPCHEM1\BNA F\DATA\BF032718\
Data File : BF103979.D
Acq On : 27 Mar 2018 19:29
Operator : JU/SJ
Sample : J2065-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
RO-CV-2536

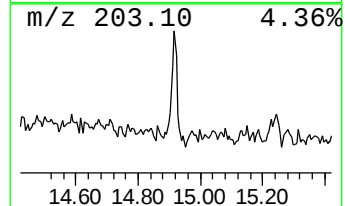
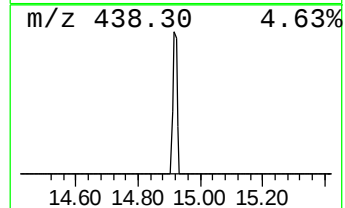
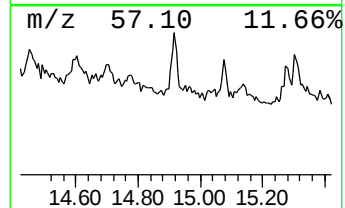
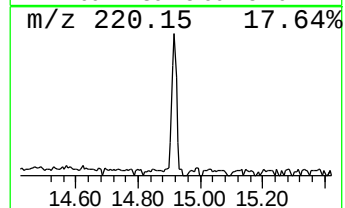
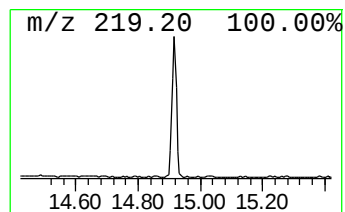
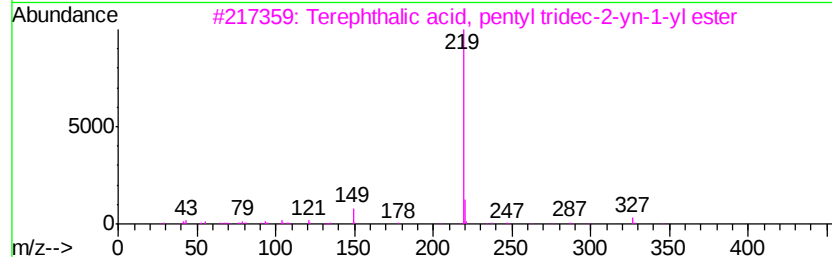
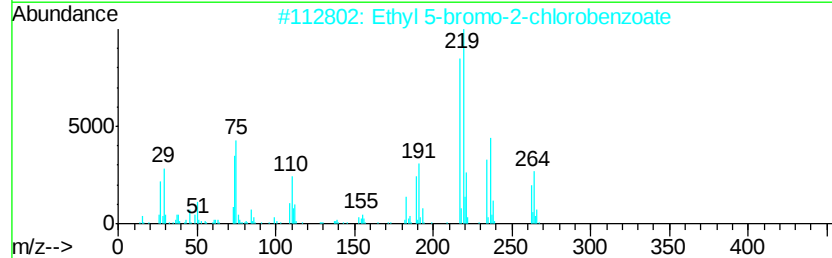
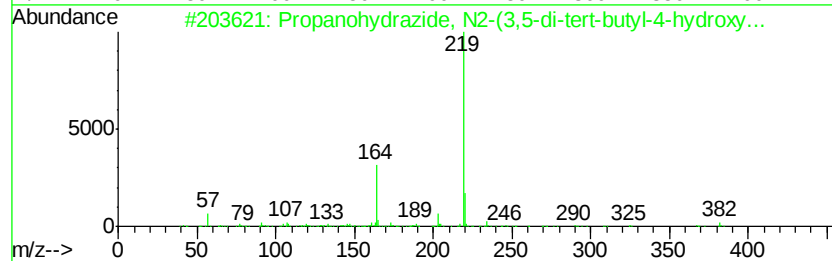
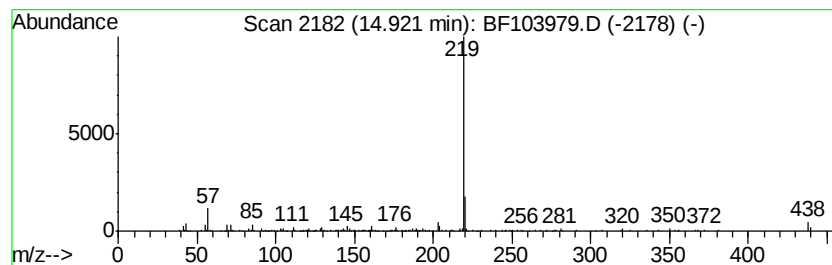
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF031518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 Propanohydrazide, N2-(3,5-d... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.92	3.88 ng	128283	Chrysene-d12	14.16

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propanohydrazide, N2-(3,5-di-ter...	382	C24H34N2O2	304870-86-8	64
2			Ethyl 5-bromo-2-chlorobenzoate	262	C9H8BrClO2	1000343-08-1	59
3			Terephthalic acid, pentyl tridec...	414	C26H38O4	1000323-92-6	59
4			Terephthalic acid, 2-isopropylph...	354	C22H26O4	1000323-58-2	50
5			4,4'-Ethylenebis(2,6-di-tert-but...	438	C30H46O2	001516-94-5	50



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 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RO-CV-2536

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-methoxy...	2.32	3.8	ng	142174	1	6.97	746554	20.0
2-Pentanone, 4-hy...	5.22	3.2	ng	117878	1	6.97	746554	20.0
unknown6.75	6.75	79.9	ng	2983820	1	6.97	746554	20.0
unknown9.20	9.20	2.5	ng	171420	3	10.02	1350880	20.0
unknown9.24	9.24	2.2	ng	149063	3	10.02	1350880	20.0
Decahydro-4,4,8,9...	9.40	5.6	ng	377734	3	10.02	1350880	20.0
unknown9.63	9.63	2.1	ng	144205	3	10.02	1350880	20.0
Naphthalene, deca...	9.66	5.5	ng	374641	3	10.02	1350880	20.0
Decahydro-8a-ethy...	9.87	23.5	ng	1589380	3	10.02	1350880	20.0
unknown9.96	9.96	9.7	ng	657373	3	10.02	1350880	20.0
1H-Indene, 2,3-di...	10.15	11.4	ng	771394	3	10.02	1350880	20.0
Naphthalene, 1,6,...	10.26	9.5	ng	643808	3	10.02	1350880	20.0
2,11-Dioxabicyclo...	10.43	30.9	ng	2087930	3	10.02	1350880	20.0
Propanohydrazide,...	14.92	3.9	ng	128283	5	14.16	660945	20.0