

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121319\
 Data File : BF118190.D
 Acq On : 13 Dec 2019 17:07
 Operator : JU
 Sample : K6235-03
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-03

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:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF120619.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	2.111	3	4	23	rVB	824549	770991	18.09%	2.490%
2	3.210	186	191	202	rVB2	29127	56736	1.33%	0.183%
3	3.605	251	258	266	rVB	41275	65905	1.55%	0.213%
4	4.981	487	492	497	rBV	81147	83292	1.95%	0.269%
5	5.069	503	507	514	rBV	134931	169882	3.99%	0.549%
6	5.481	573	577	590	rBV	1938970	1780550	41.79%	5.751%
7	5.781	624	628	632	rVB	78761	71138	1.67%	0.230%
8	6.087	675	680	683	rBV	113384	97427	2.29%	0.315%
9	6.246	704	707	712	rBV	46910	51313	1.20%	0.166%
10	6.498	746	750	758	rBV	1150009	1196728	28.09%	3.866%
11	6.640	769	774	777	rBV	3508025	3081381	72.32%	9.953%
12	6.863	809	812	816	rBV	761219	666363	15.64%	2.152%
13	7.022	835	839	843	rBV	2863430	2468163	57.93%	7.972%
14	7.434	902	909	912	rBV	1912069	1901776	44.63%	6.143%
15	7.463	912	914	919	rVB	198136	157054	3.69%	0.507%
16	7.898	981	988	991	rBV	514962	444586	10.43%	1.436%
17	8.151	1027	1031	1038	rBV	935019	829242	19.46%	2.679%
18	9.234	1210	1215	1218	rBV	4033881	3686035	86.51%	11.906%
19	9.622	1278	1281	1286	rBV	91219	83331	1.96%	0.269%
20	9.916	1327	1331	1341	rVB	1055336	989901	23.23%	3.197%
21	10.216	1377	1382	1386	rBV4	24937	49416	1.16%	0.160%
22	10.710	1461	1466	1469	rBV	3070564	2755129	64.66%	8.899%
23	11.139	1534	1539	1547	rVB	94280	110408	2.59%	0.357%
24	11.404	1581	1584	1592	rBV	1117055	1014280	23.80%	3.276%
25	11.933	1669	1674	1681	rBV	871069	864840	20.30%	2.794%
26	11.992	1681	1684	1690	rVB2	56403	75610	1.77%	0.244%
27	12.716	1804	1807	1814	rBV	547496	586206	13.76%	1.894%
28	12.998	1851	1855	1858	rBV	4464112	4260963	100.00%	13.763%
29	13.886	2003	2006	2015	rBV	247334	278432	6.53%	0.899%
30	14.057	2031	2035	2044	rBV	826258	789342	18.52%	2.550%
31	14.210	2058	2061	2066	rVB	69594	65579	1.54%	0.212%
32	14.516	2110	2113	2121	rBV2	99282	96722	2.27%	0.312%
33	14.827	2162	2166	2172	rBV	117071	117269	2.75%	0.379%
34	15.174	2221	2225	2231	rVB	108895	128434	3.01%	0.415%

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Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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ClientSampleId :
MW-03

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:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF120619.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	15.551	2284	2289	2300	rVB	616796	892862	20.95%	2.884%
36	16.010	2362	2367	2373	rBV2	87996	145183	3.41%	0.469%
37	16.527	2450	2455	2460	rVB2	44745	76388	1.79%	0.247%

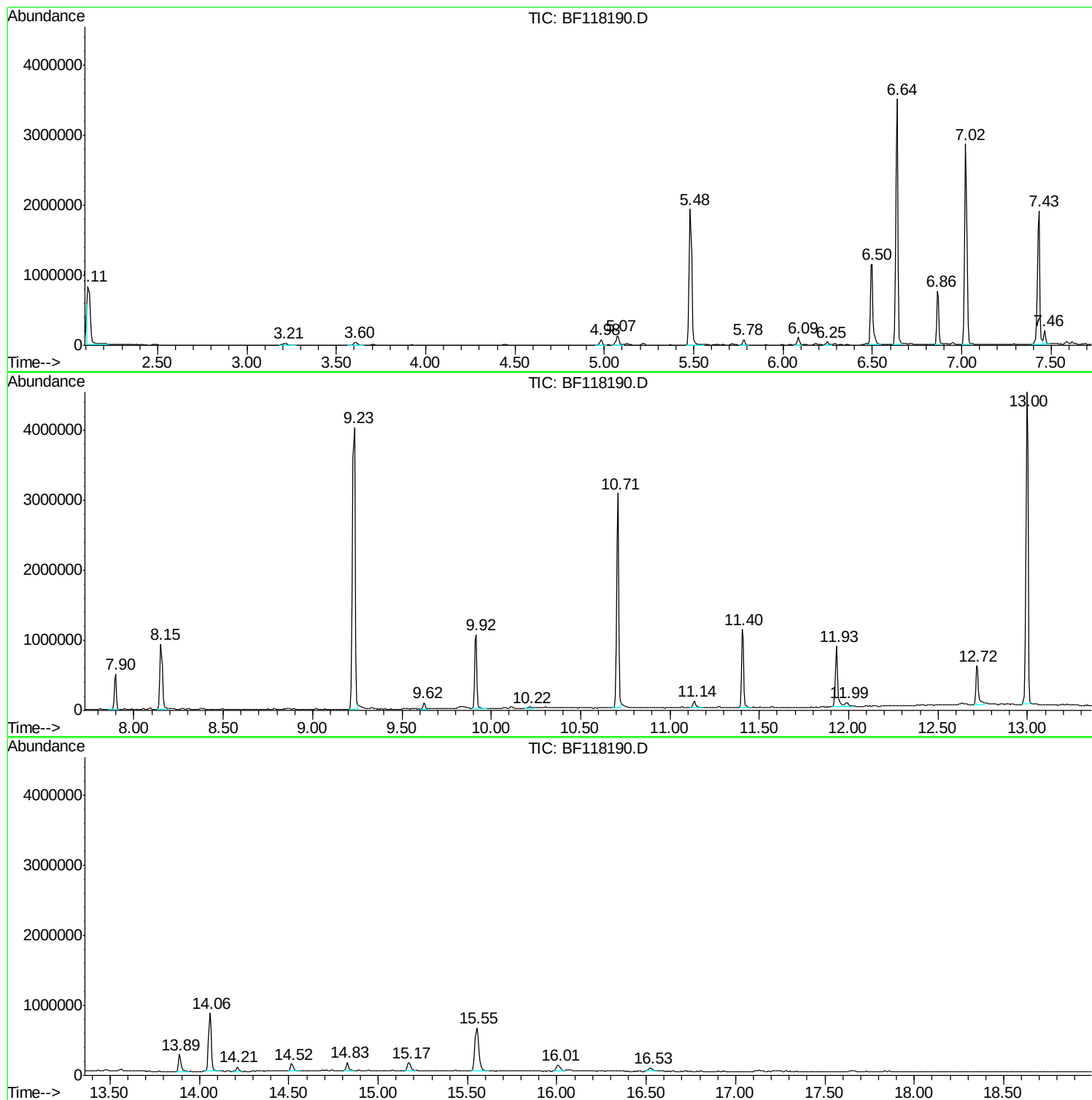
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ClientSampled :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF120619.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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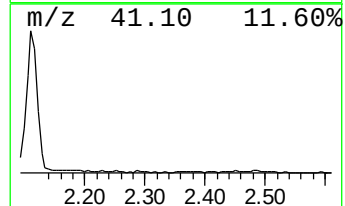
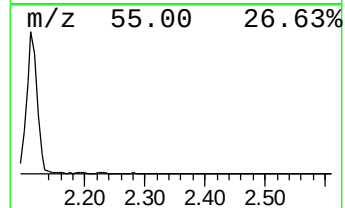
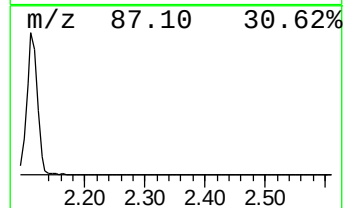
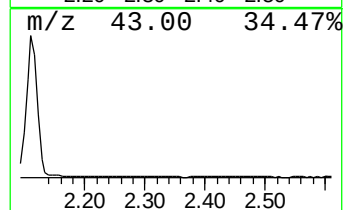
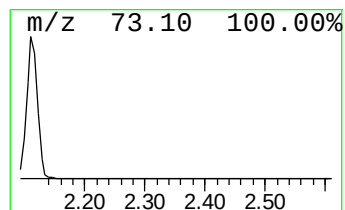
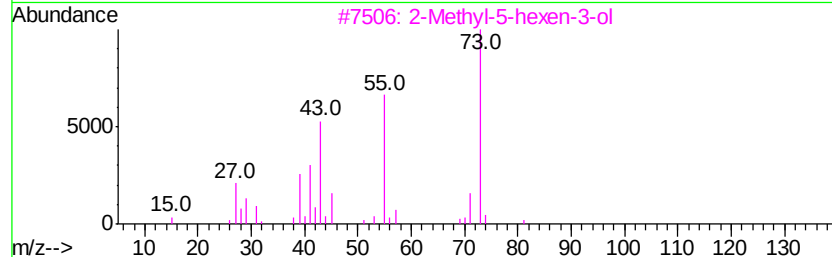
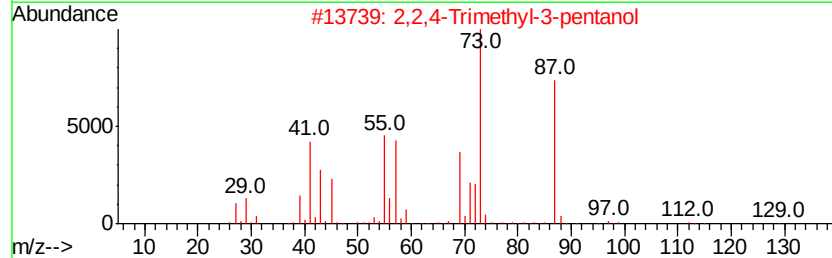
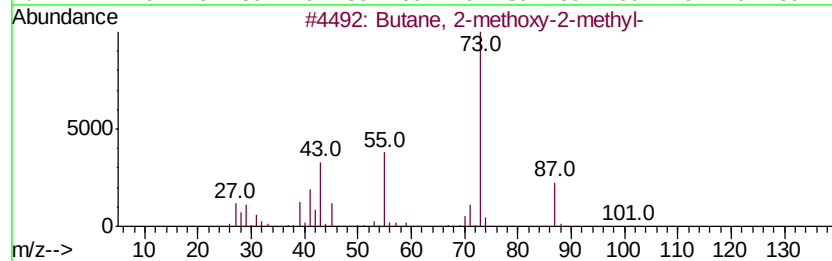
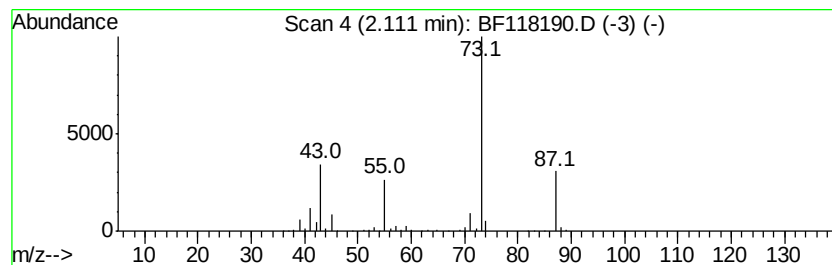
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF120619.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.11	23.14 ng	770991	1,4-Dichlorobenzene-d4	6.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
3		2-Methyl-5-hexen-3-ol	114	C7H14O	032815-70-6	23
4		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
5		Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	12



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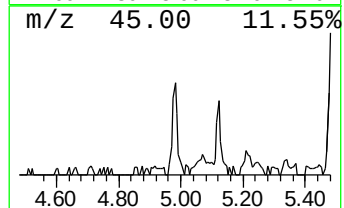
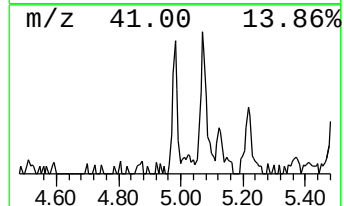
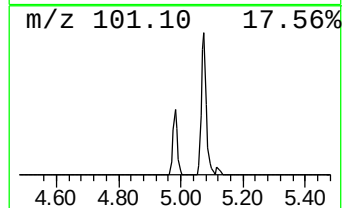
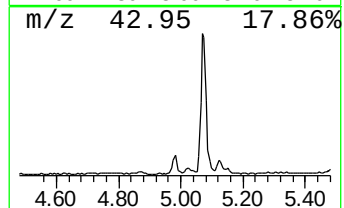
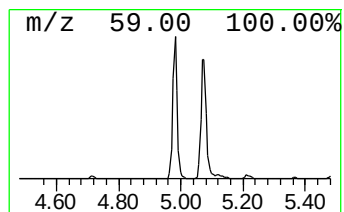
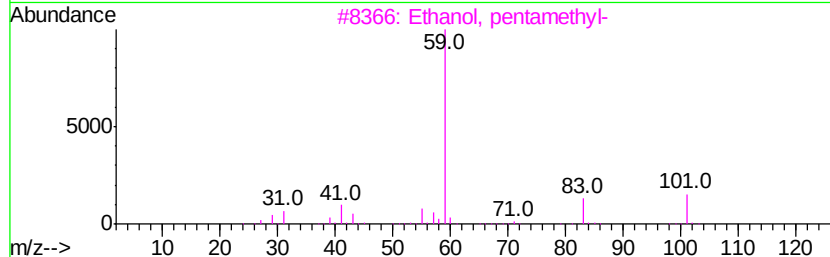
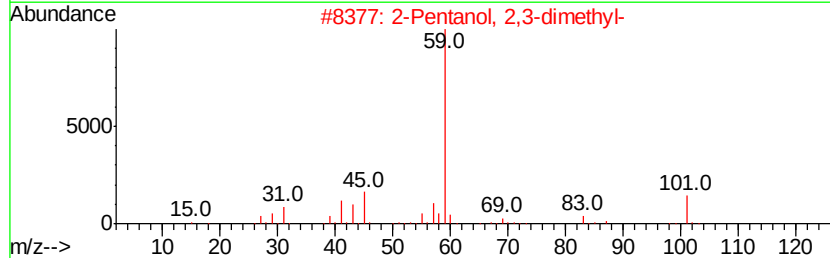
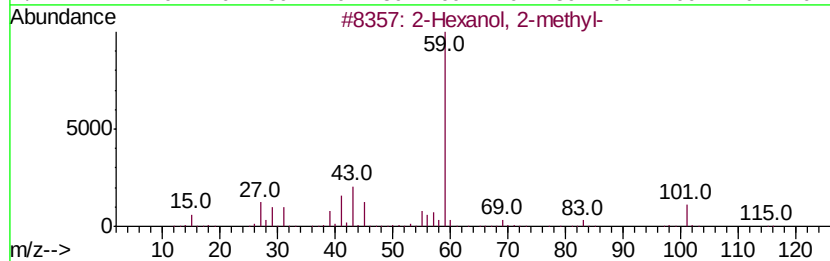
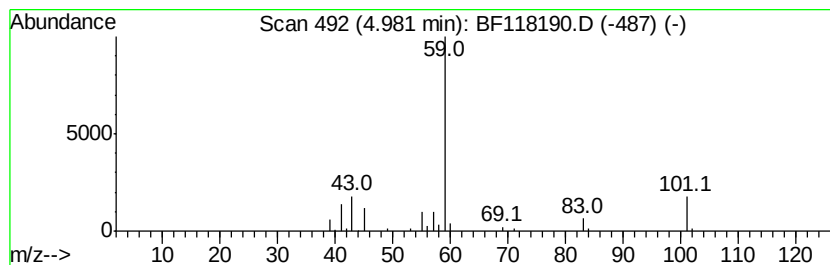
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 2-Hexanol, 2-methyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.98	2.50 ng	83292	1,4-Dichlorobenzene-d4	6.86

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	78
2			2-Pentanol, 2,3-dimethyl-	116	C7H16O	004911-70-0	74
3			Ethanol, pentamethyl-	116	C7H16O	000594-83-2	56
4			3-Heptanol, 4-methyl-	130	C8H18O	014979-39-6	40
5			2-Pentanol, 2,4-dimethyl-	116	C7H16O	000625-06-9	40



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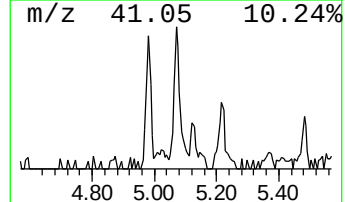
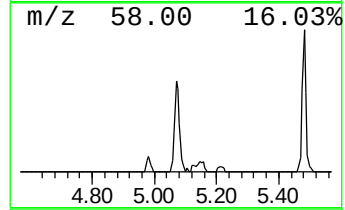
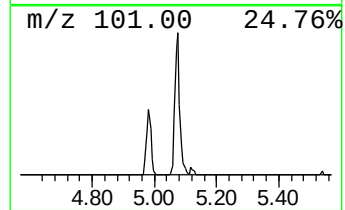
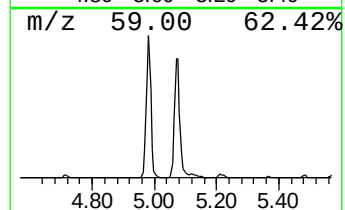
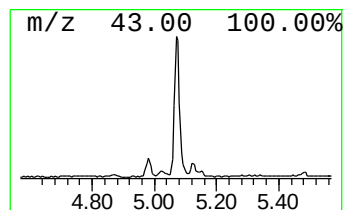
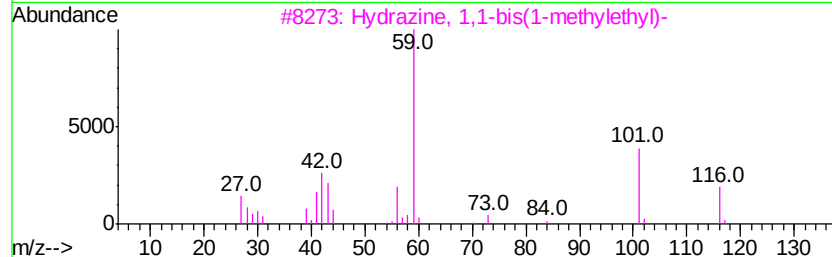
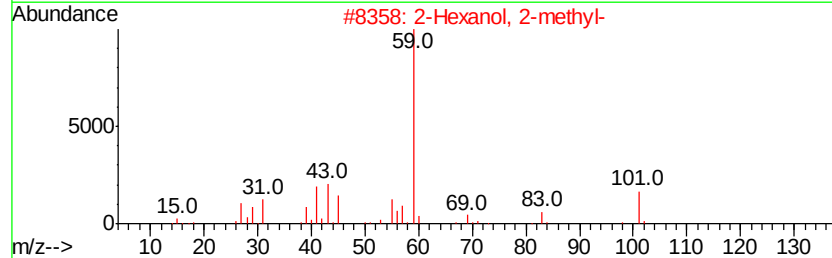
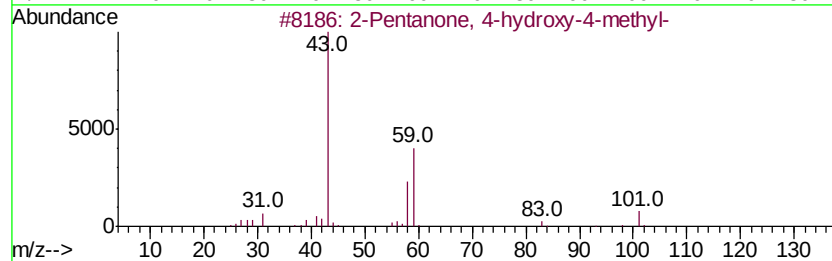
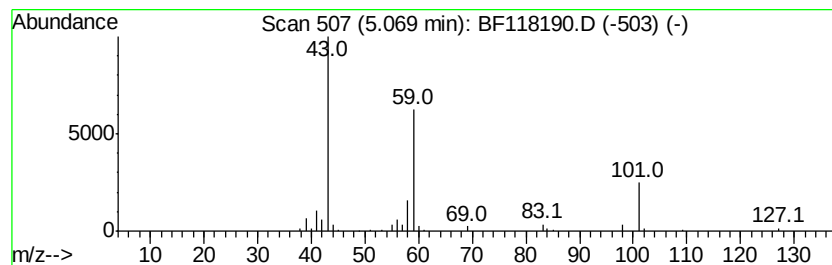
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.07	5.10 ng	169882	1,4-Dichlorobenzene-d4	6.86

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	40
3			Hydrazine, 1,1-bis(1-methylethyl)-	116	C6H16N2	000921-14-2	38
4			Acetic acid, cyano-, 1,1-dimethyl-	141	C7H11NO2	001116-98-9	25
5			Methyl 2,5,8,11-tetraoxatridecan-	236	C10H20O6	1000366-78-2	23



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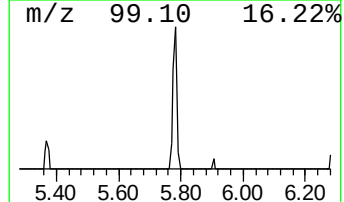
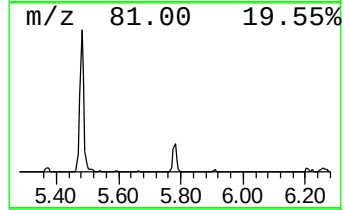
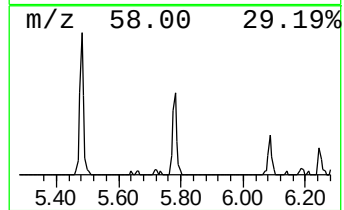
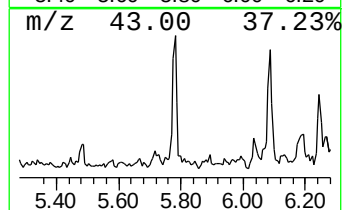
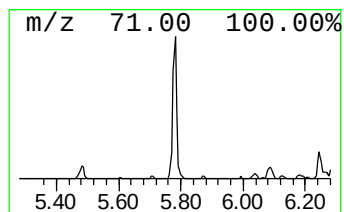
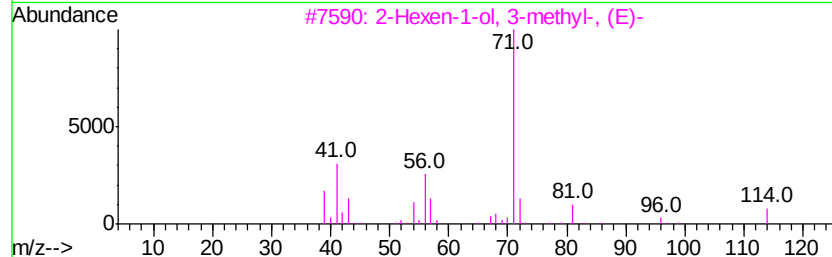
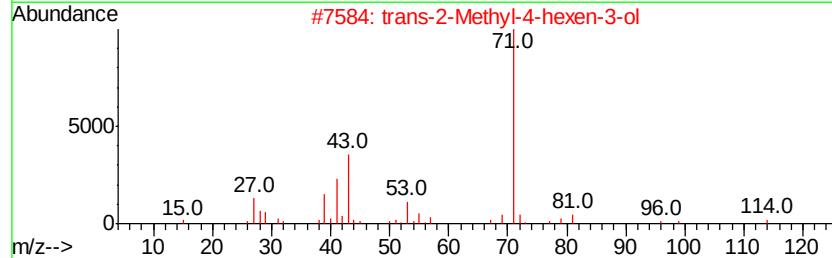
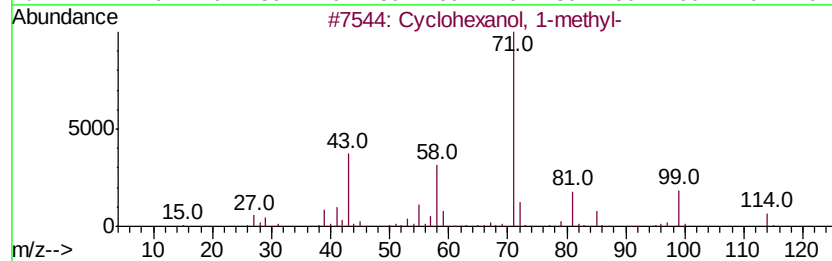
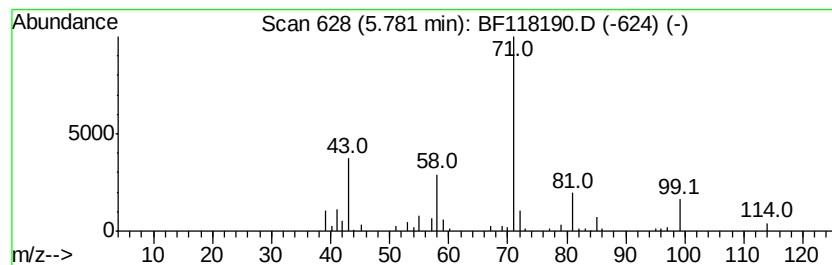
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 Cyclohexanol, 1-methyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.78	2.14 ng	71138	1,4-Dichlorobenzene-d4	6.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexanol, 1-methyl-	114	C7H14O	000590-67-0	91
2		trans-2-Methyl-4-hexen-3-ol	114	C7H14O	096346-76-8	50
3		2-Hexen-1-ol, 3-methyl-, (E)-	114	C7H14O	030801-96-8	40
4		1,2-Dimethylcyclopentanol	114	C7H14O	019550-45-9	37
5		2,4-Dimethylcyclopentanol	114	C7H14O	089794-28-5	36



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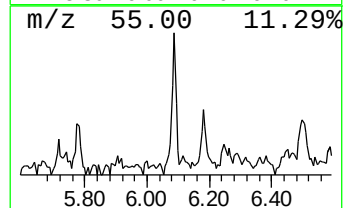
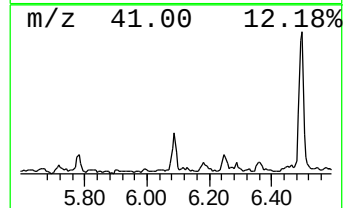
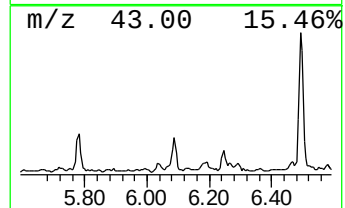
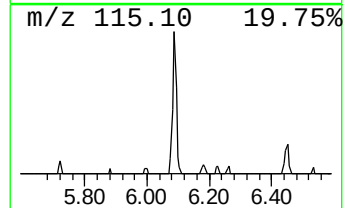
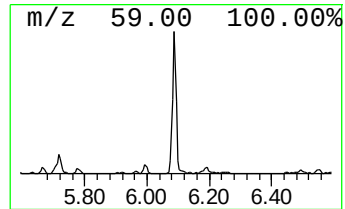
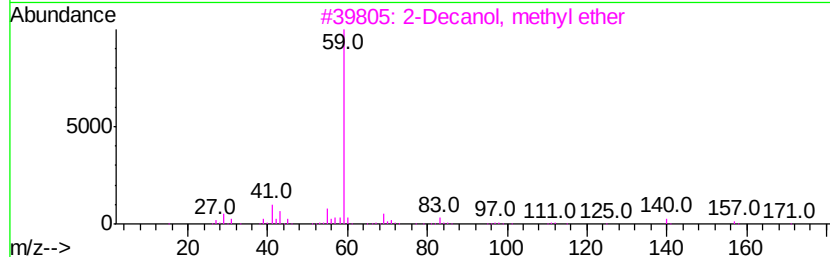
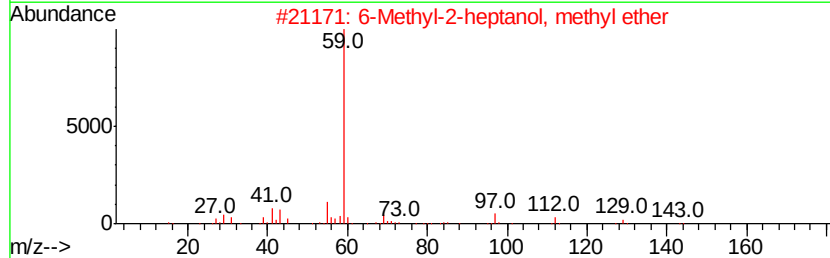
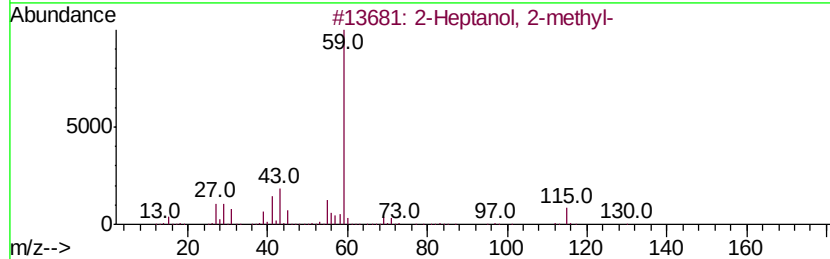
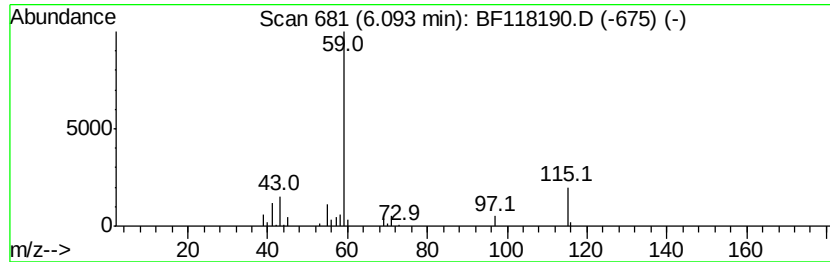
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 2-Heptanol, 2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.09	2.92 ng	97427	1,4-Dichlorobenzene-d4	6.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Heptanol, 2-methyl-	130	C8H18O	000625-25-2	78
2		6-Methyl-2-heptanol, methyl ether	144	C9H20O	1000365-52-0	72
3		2-Decanol, methyl ether	172	C11H24O	1000333-83-9	64
4		2-Hexanol, 2,3-dimethyl-	130	C8H18O	019550-03-9	56
5		Butane, 2-methoxy-3-methyl-	102	C6H14O	062016-49-3	45



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Sample : K6235-03
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-03

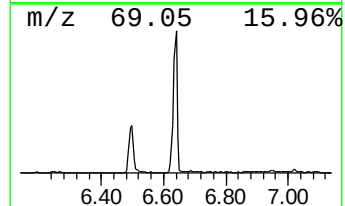
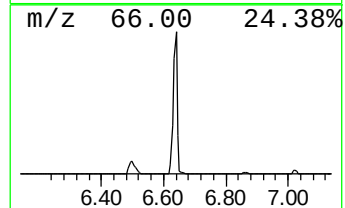
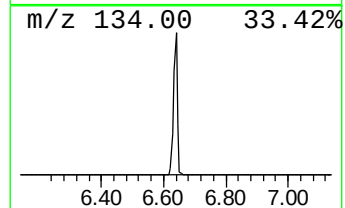
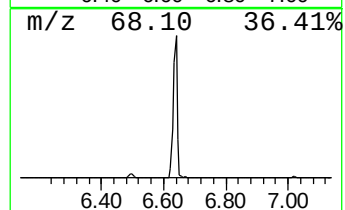
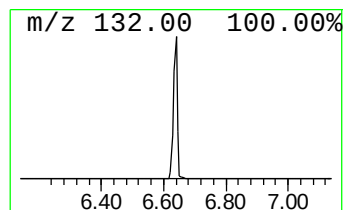
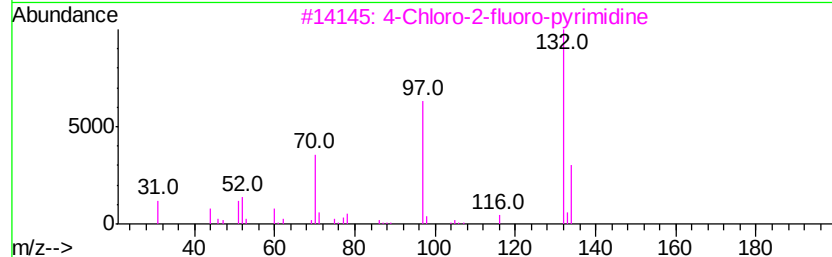
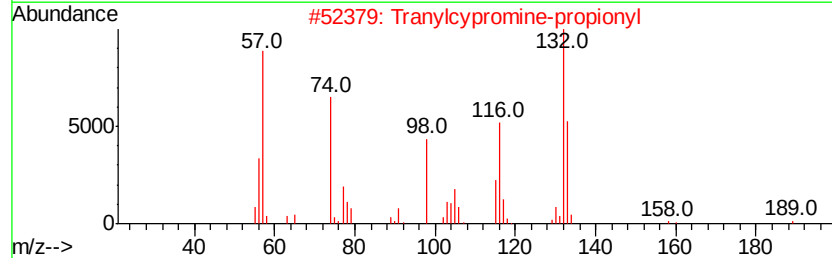
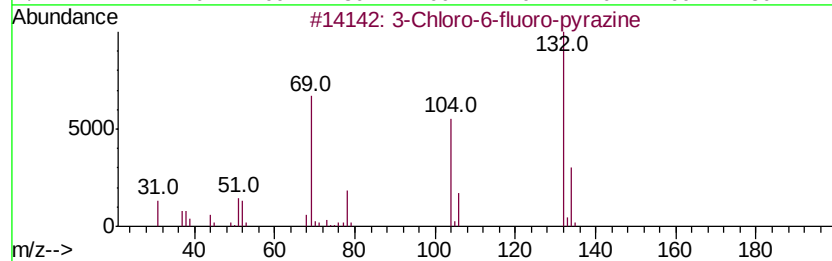
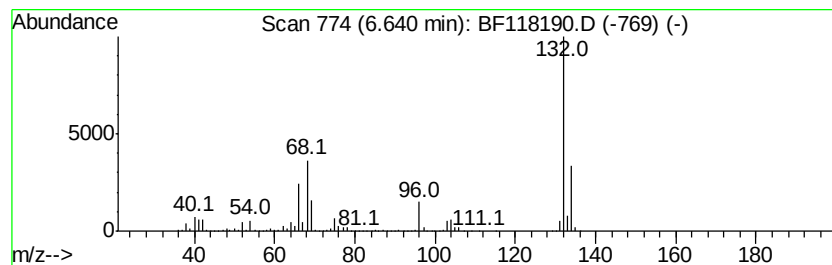
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 unknown6.64 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.64	92.48 ng	3081380	1,4-Dichlorobenzene-d4	6.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	16
3		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9
4		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121319\
Data File : BF118190.D
Acq On : 13 Dec 2019 17:07
Operator : JU
Sample : K6235-03
Misc :
ALS Vial : 14 Sample Multiplier: 1

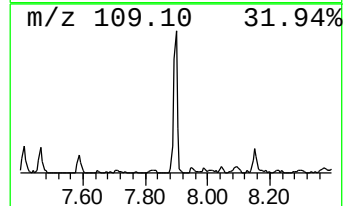
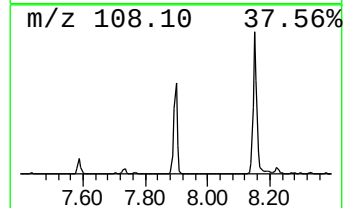
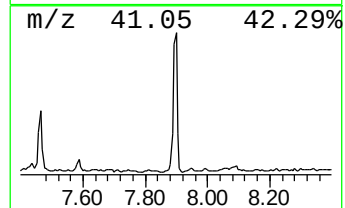
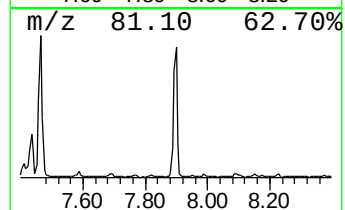
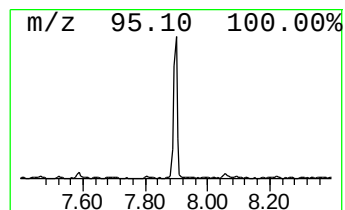
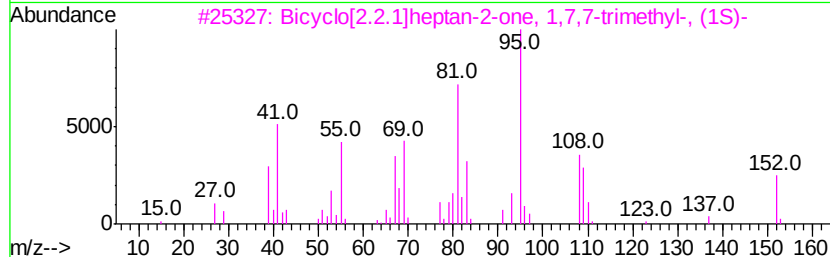
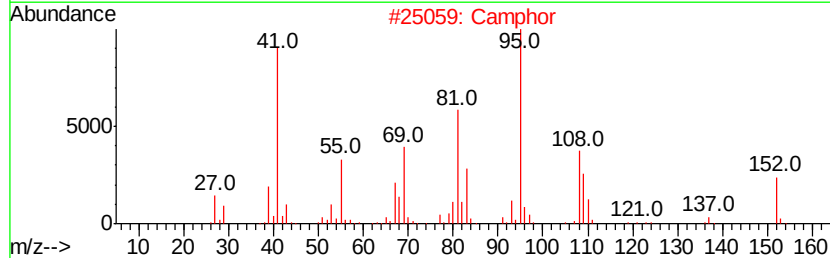
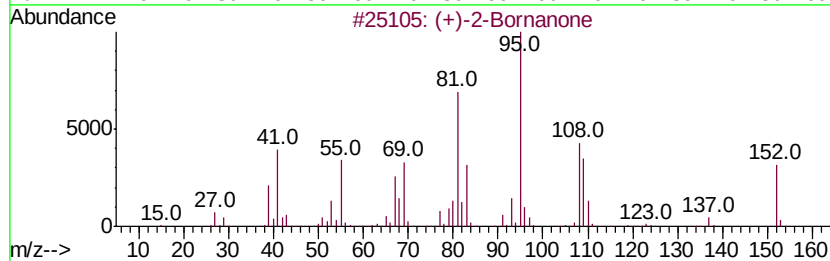
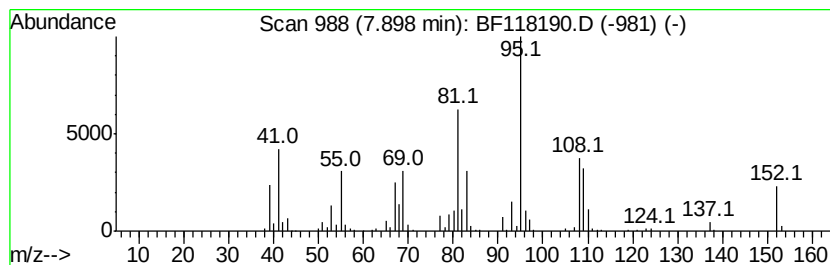
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 (+)-2-Bornanone Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
7.90	10.72 ng	444586	Napthalene-d8	8.15	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(+)-2-Bornanone	152	C10H16O	000464-49-3	98
2	Camphor	152	C10H16O	000076-22-2	97
3	Bicyclo[2.2.1]heptan-2-one, 1,7,...	152	C10H16O	000464-48-2	96
4	2-Butanone, 4-(5-methyl-2-furanyl)-	152	C9H12O2	013679-56-6	49
5	2-Furancarboxylic acid, 2-propen...	152	C8H8O3	004208-49-5	43



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121319\
Data File : BF118190.D
Acq On : 13 Dec 2019 17:07
Operator : JU
Sample : K6235-03
Misc :
ALS Vial : 14 Sample Multiplier: 1

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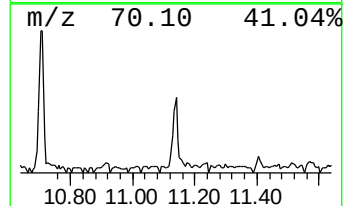
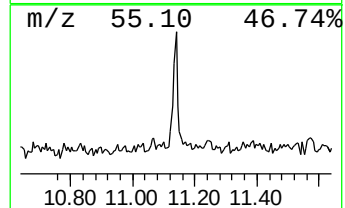
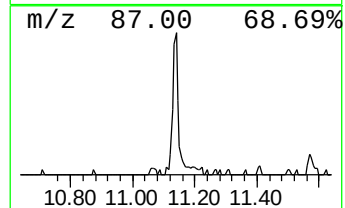
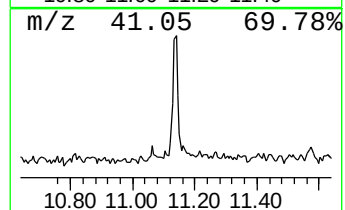
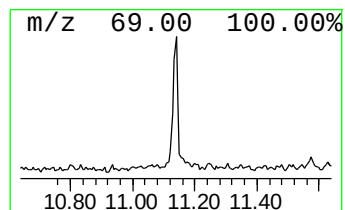
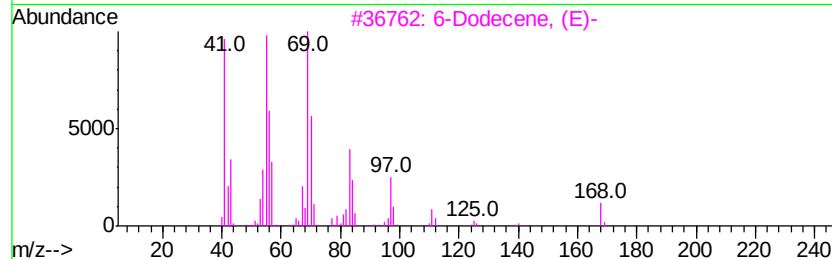
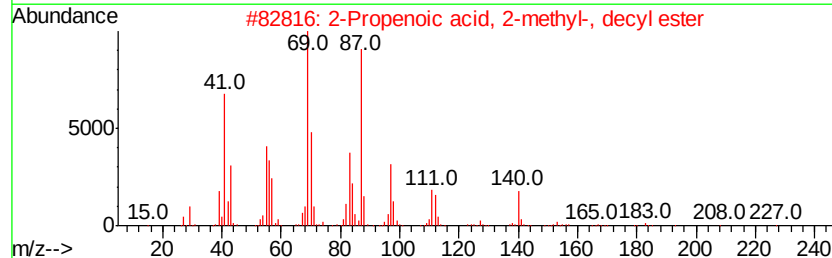
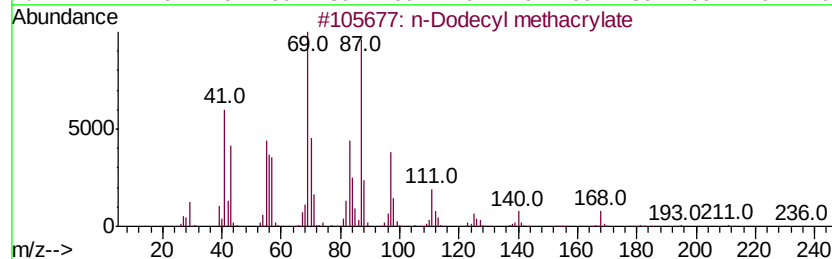
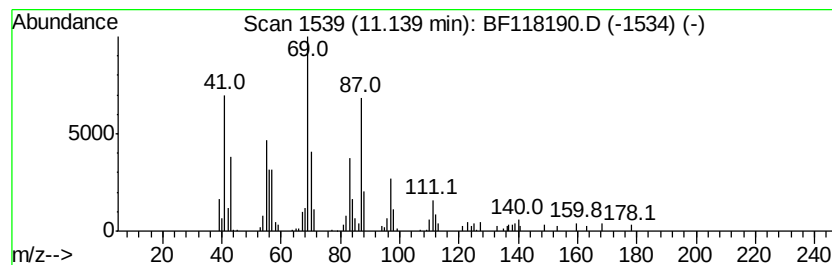
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 n-Dodecyl methacrylate Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.14	2.18 ng	110408	Phenanthrene-d10	11.40

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Dodecyl methacrylate	254	C16H30O2	000142-90-5	91
2	2-Propenoic acid, 2-methyl-, dec...	226	C14H26O2	003179-47-3	91
3	6-Dodecene, (E)-	168	C12H24	007206-17-9	59
4	Cyclopropanecarboxylic acid, non...	212	C13H24O2	060128-06-5	59
5	1-Pentene, 3-ethyl-	98	C7H14	004038-04-4	53



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121319\
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Sample : K6235-03
Misc :
ALS Vial : 14 Sample Multiplier: 1

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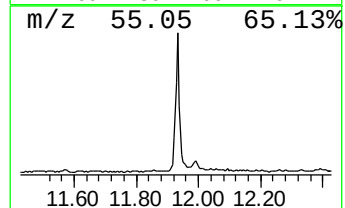
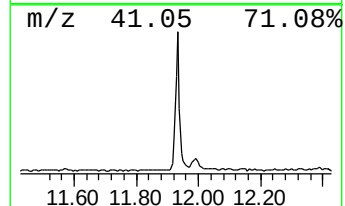
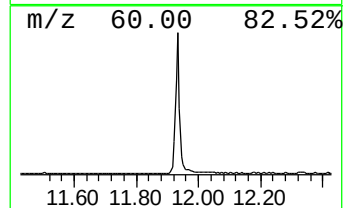
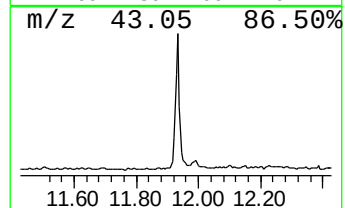
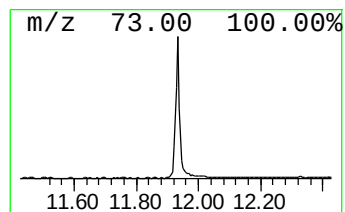
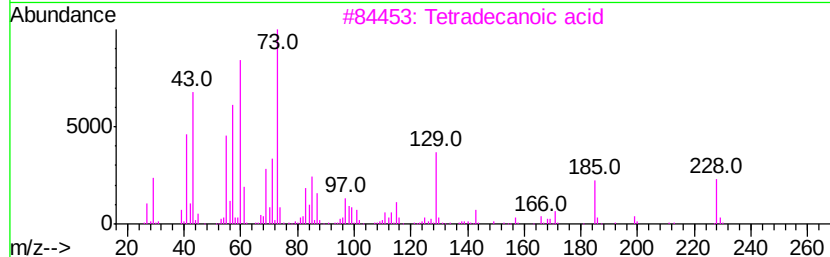
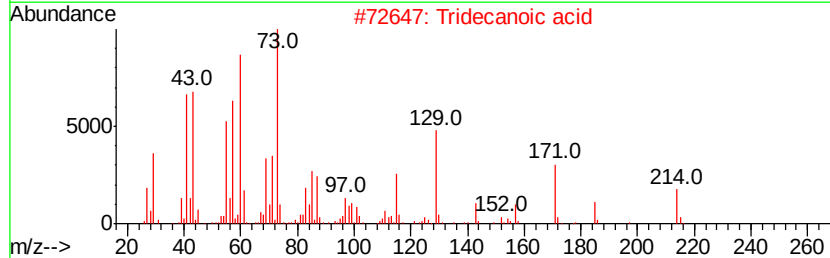
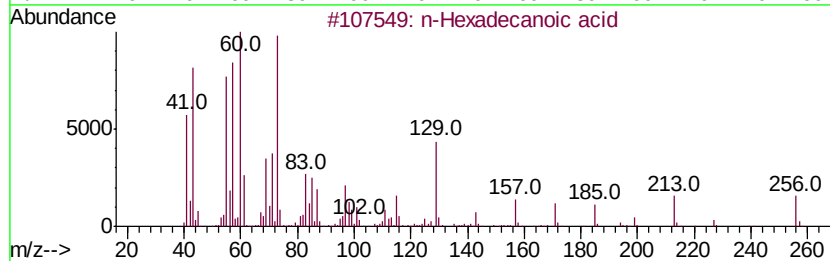
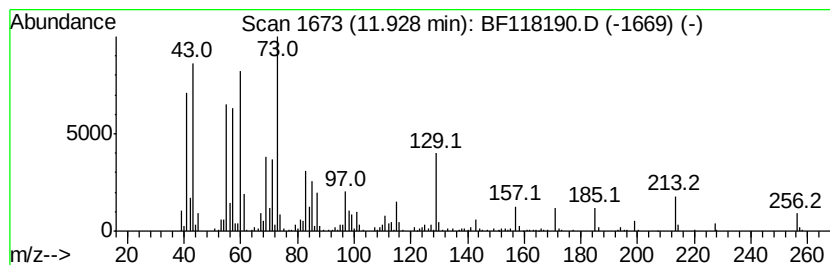
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.93	17.05 ng	864840	Phenanthrene-d10	11.40

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		Tridecanoic acid	214	C13H26O2	000638-53-9	94
3		Tetradecanoic acid	228	C14H28O2	000544-63-8	93
4		Pentadecanoic acid	242	C15H30O2	001002-84-2	91
5		Decanoic acid, silver(1+) salt	278	C10H19AgO2	013126-67-5	59



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121319\
Data File : BF118190.D
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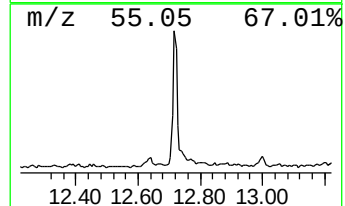
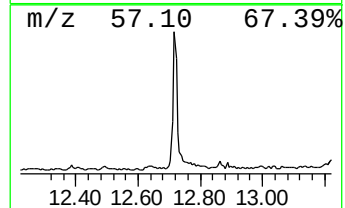
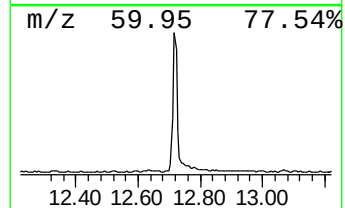
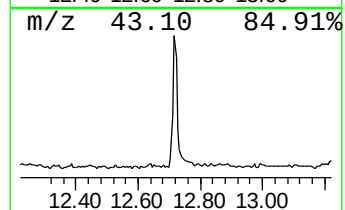
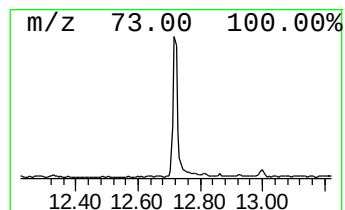
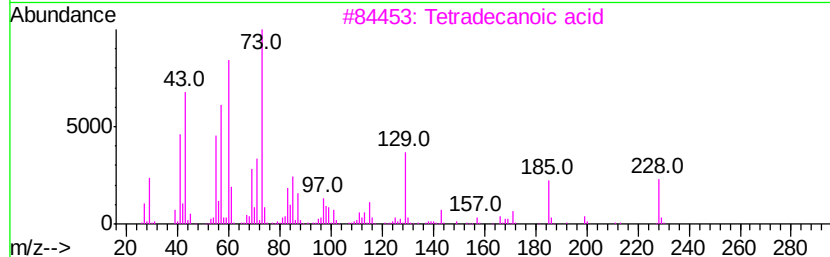
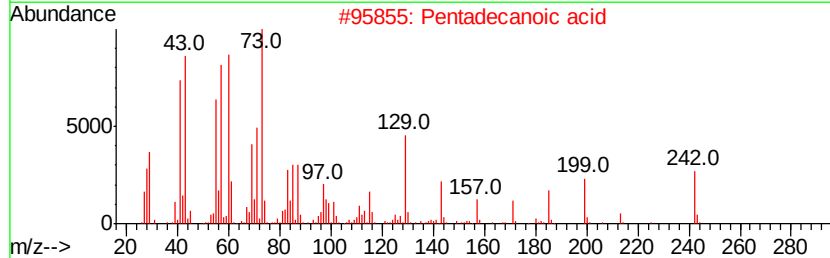
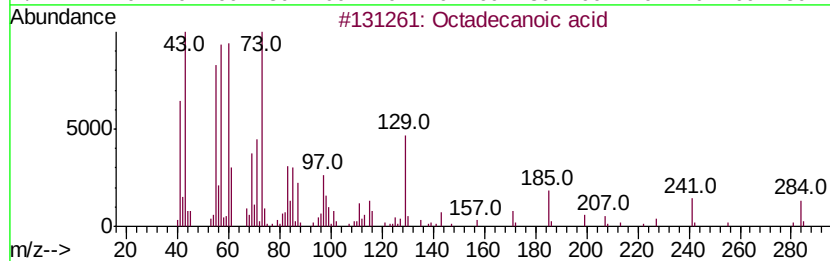
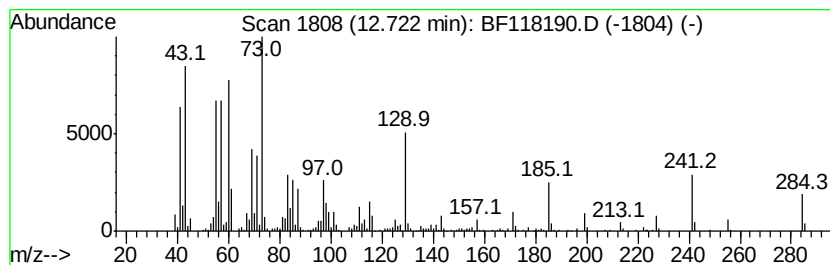
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Octadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.72	11.56 ng	586206	Phenanthrene-d10	11.40

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2		Pentadecanoic acid	242	C15H30O2	001002-84-2	93
3		Tetradecanoic acid	228	C14H28O2	000544-63-8	74
4		n-Decanoic acid	172	C10H20O2	000334-48-5	70
5		Tridecanoic acid	214	C13H26O2	000638-53-9	68



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121319\
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ClientSampleId :
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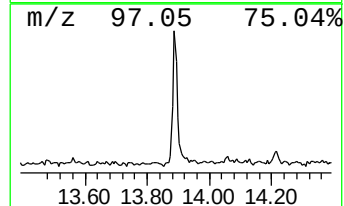
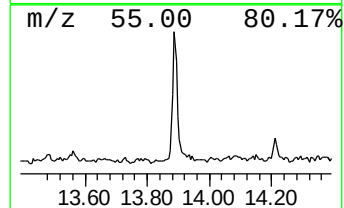
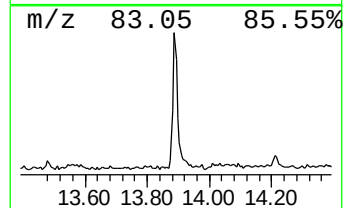
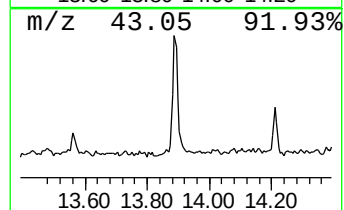
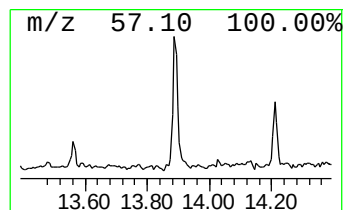
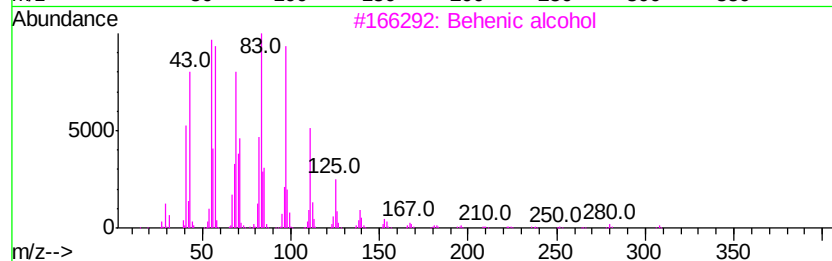
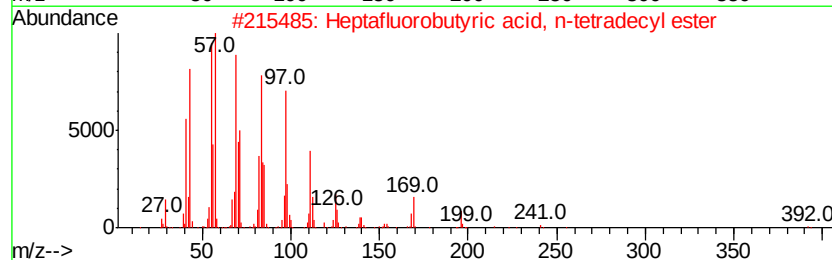
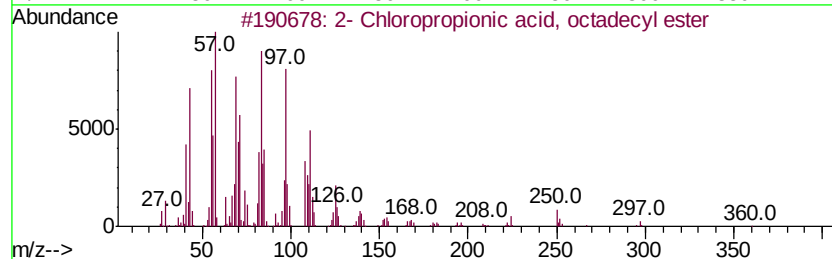
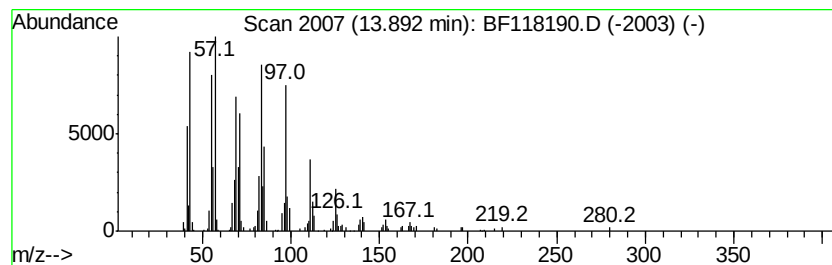
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 2- Chloropropionic acid, oc... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.89	7.05 ng	278432	Chrysene-d12	14.06

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2- Chloropropionic acid, octadec...	360	C21H41ClO2	088104-31-8	94
2			Heptafluorobutyric acid, n-tetra...	410	C18H29F7O2	007365-36-8	94
3			Behenic alcohol	326	C22H46O	000661-19-8	94
4			n-Tetracosanol-1	354	C24H50O	000506-51-4	94
5			Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	94



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121319\
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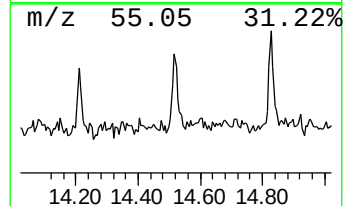
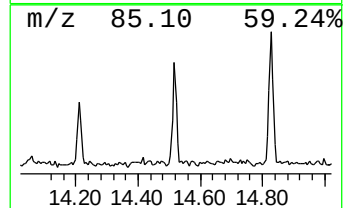
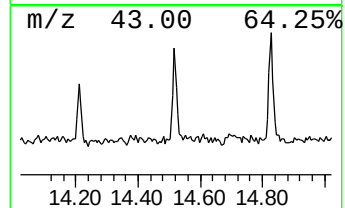
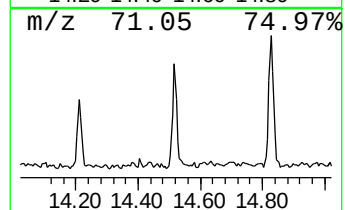
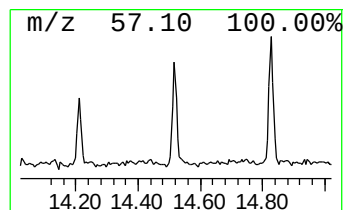
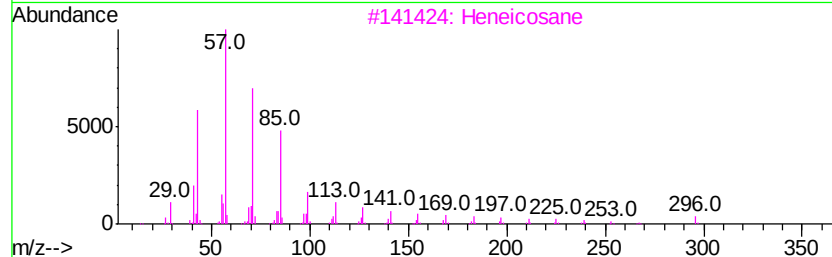
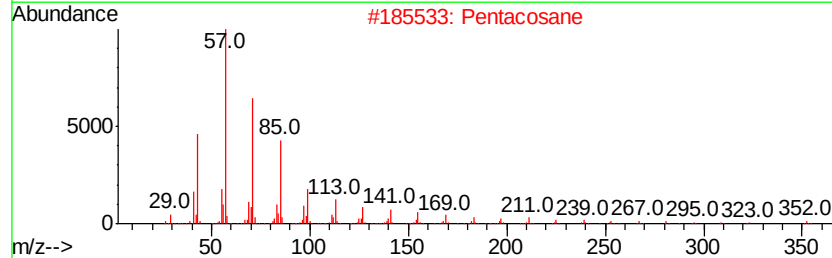
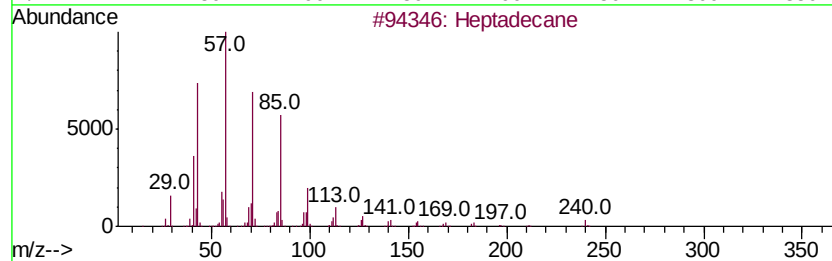
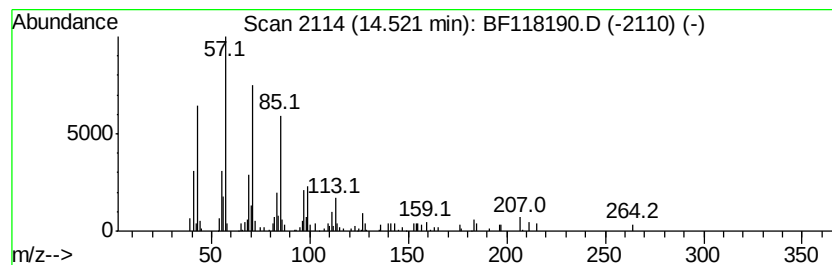
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 Heptadecane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.52	2.45 ng	96722	Chrysene-d12	14.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane	240	C17H36	000629-78-7	91
2		Pentacosane	352	C25H52	000629-99-2	90
3		Heneicosane	296	C21H44	000629-94-7	90
4		Octacosane	394	C28H58	000630-02-4	90
5		Triacontane	422	C30H62	000638-68-6	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121319\
Data File : BF118190.D
Acq On : 13 Dec 2019 17:07
Operator : JU
Sample : K6235-03
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-03

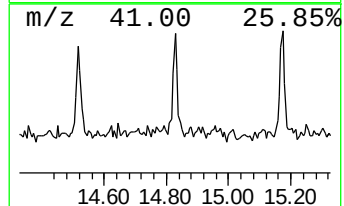
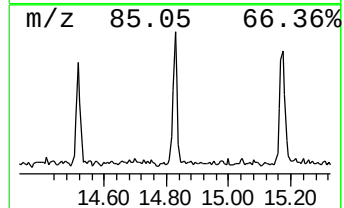
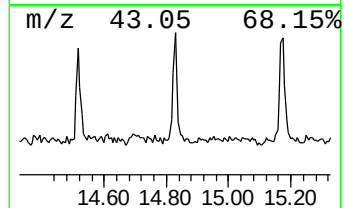
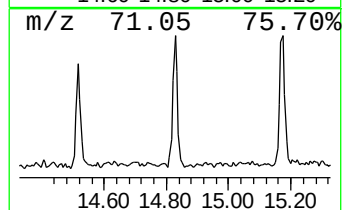
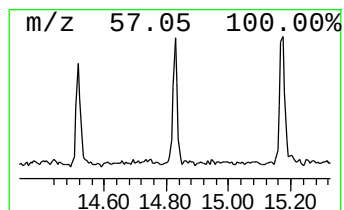
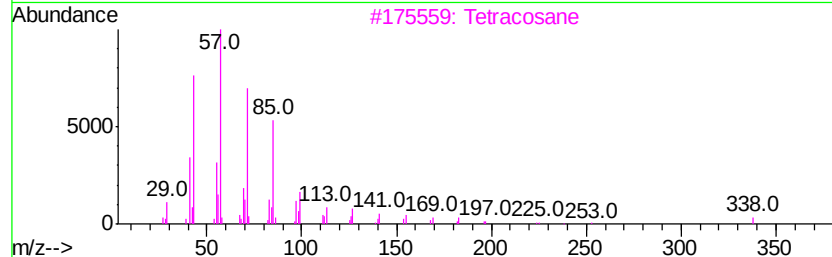
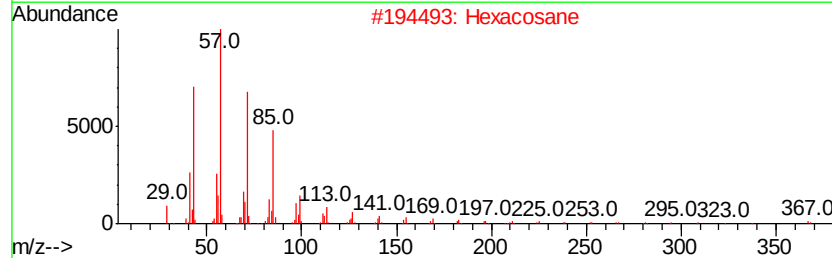
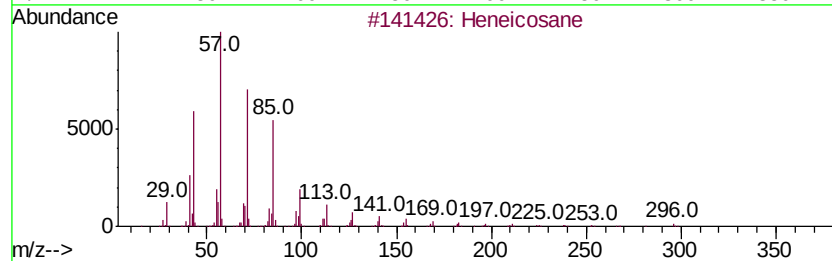
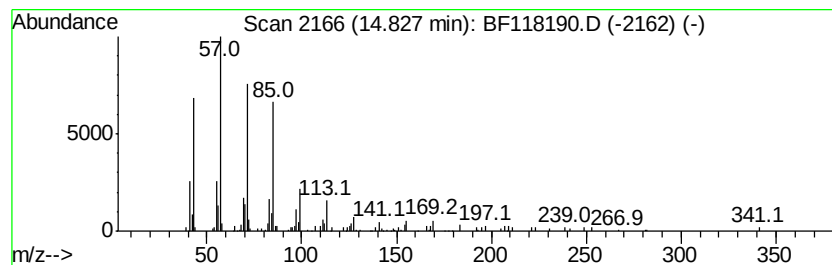
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Heneicosane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.83	2.63 ng	117269	Perylene-d12	15.55

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	91
2		Hexacosane	366	C26H54	000630-01-3	90
3		Tetracosane	338	C24H50	000646-31-1	90
4		Octacosane	394	C28H58	000630-02-4	90
5		triacontane	422	C30H62	000638-68-6	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121319\
Data File : BF118190.D
Acq On : 13 Dec 2019 17:07
Operator : JU
Sample : K6235-03
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-03

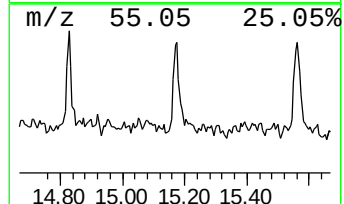
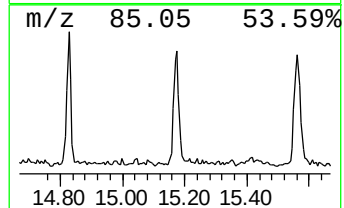
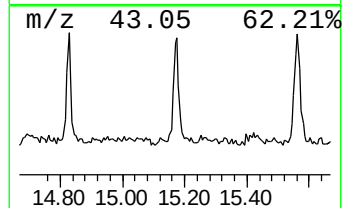
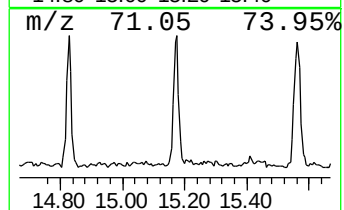
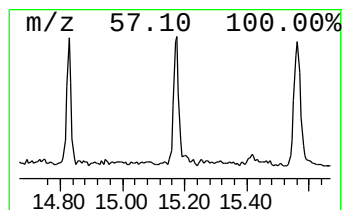
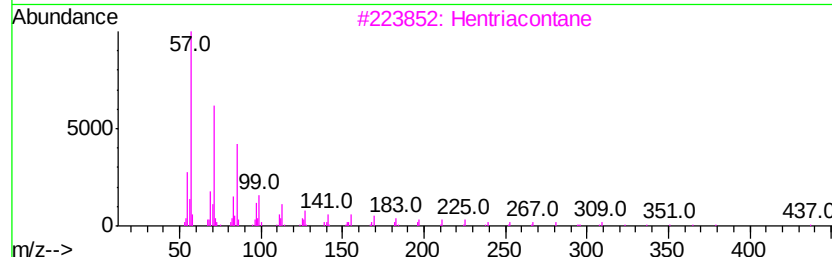
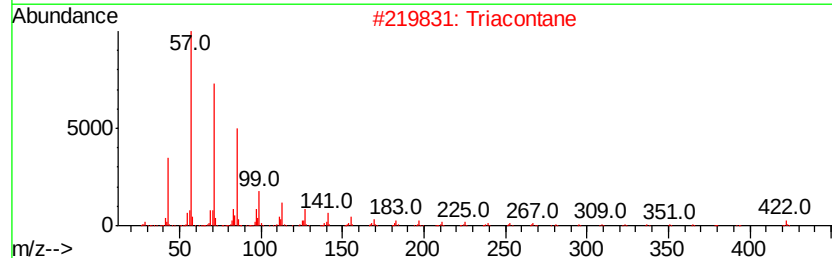
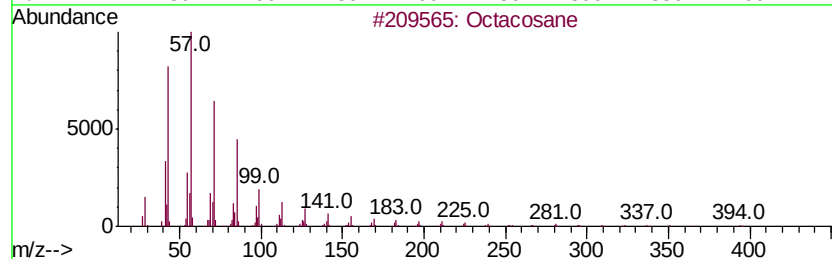
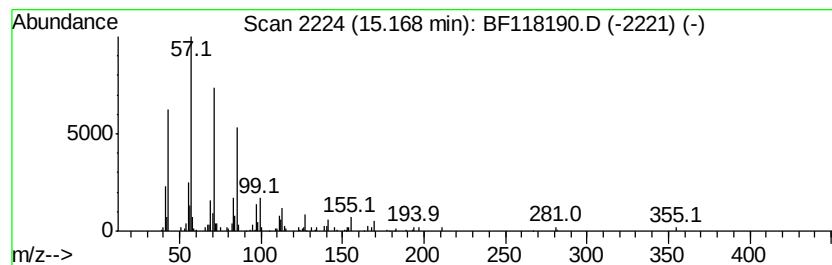
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Octacosane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.17	2.88 ng	128434	Perylene-d12	15.55

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octacosane	394	C28H58	000630-02-4	91
2		triacontane	422	C30H62	000638-68-6	91
3		Hentriacontane	437	C31H64	000630-04-6	91
4		Hexacosane	366	C26H54	000630-01-3	91
5		Heneicosane	296	C21H44	000629-94-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121319\
Data File : BF118190.D
Acq On : 13 Dec 2019 17:07
Operator : JU
Sample : K6235-03
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-03

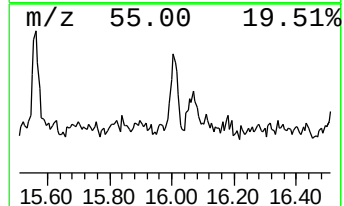
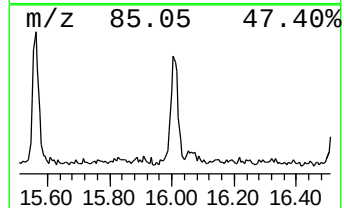
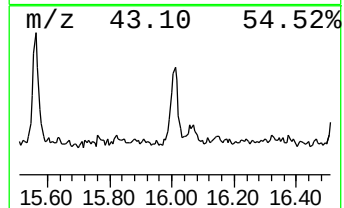
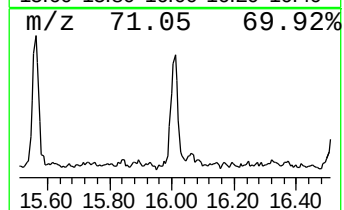
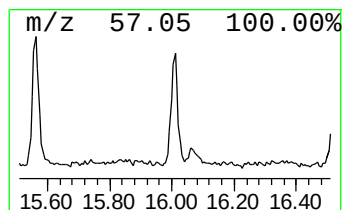
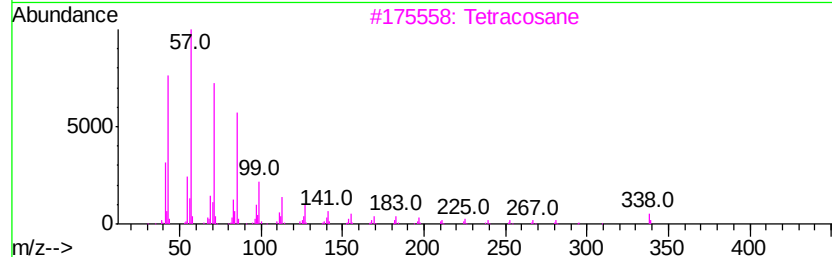
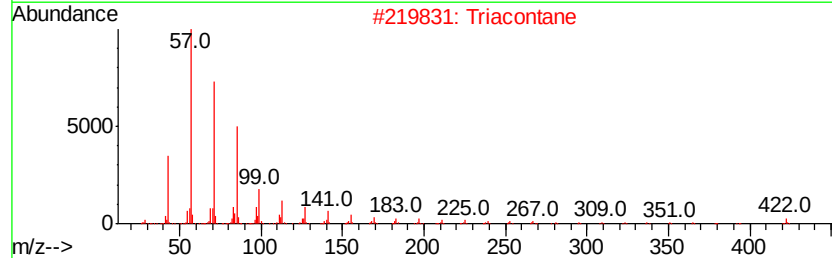
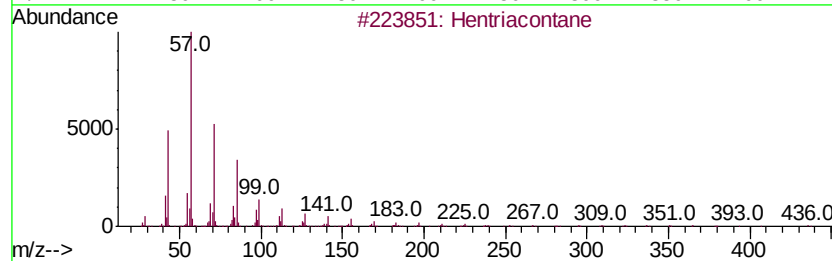
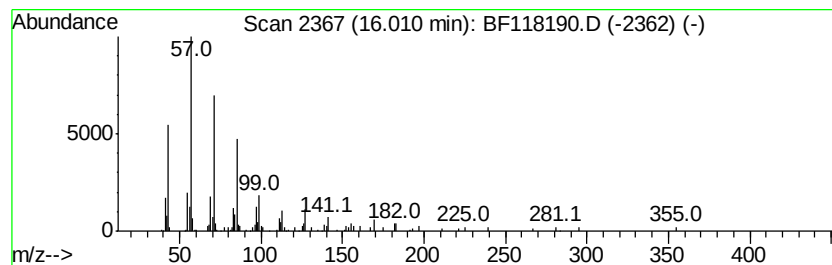
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF120619.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Hentriacontane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.01	3.25 ng	145183	Perylene-d12	15.55

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hentriacontane	437	C31H64	000630-04-6	91
2			Triacontane	422	C30H62	000638-68-6	91
3			Tetracosane	338	C24H50	000646-31-1	91
4			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
5			Heneicosane	296	C21H44	000629-94-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF121319\
 Data File : BF118190.D
 Acq On : 13 Dec 2019 17:07
 Operator : JU
 Sample : K6235-03
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-03

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF120619.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...	2.11	23.1	ng	770991	1	6.86	666363	20.0
2-Hexanol, 2-methyl-	4.98	2.5	ng	83292	1	6.86	666363	20.0
2-Pentanone, 4-hy...	5.07	5.1	ng	169882	1	6.86	666363	20.0
Cyclohexanol, 1-m...	5.78	2.1	ng	71138	1	6.86	666363	20.0
2-Heptanol, 2-met...	6.09	2.9	ng	97427	1	6.86	666363	20.0
unknown6.64	6.64	92.5	ng	3081380	1	6.86	666363	20.0
(+)-2-Bornanone	7.90	10.7	ng	444586	2	8.15	829242	20.0
n-Dodecyl methacr...	11.14	2.2	ng	110408	4	11.40	1014280	20.0
n-Hexadecanoic acid	11.93	17.1	ng	864840	4	11.40	1014280	20.0
Octadecanoic acid	12.72	11.6	ng	586206	4	11.40	1014280	20.0
2-Chloropropioni...	13.89	7.0	ng	278432	5	14.06	789342	20.0
Heptadecane	14.52	2.5	ng	96722	5	14.06	789342	20.0
Heneicosane	14.83	2.6	ng	117269	6	15.55	892862	20.0
Octacosane	15.17	2.9	ng	128434	6	15.55	892862	20.0
Hentriacontane	16.01	3.3	ng	145183	6	15.55	892862	20.0