

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF072820\
 Data File : BF121063.D
 Acq On : 28 Jul 2020 11:48
 Operator : JU/CG
 Sample : L3378-16
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 DRUM-WATER

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-BF071620.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	5.010	493	497	507	rVB	391502	382257	6.78%	1.230%
2	5.422	563	567	585	rBV	864666	986970	17.51%	3.175%
3	5.757	620	624	633	rBV	149405	139980	2.48%	0.450%
4	6.304	713	717	725	rBV	996347	1094026	19.41%	3.520%
5	6.439	736	740	743	rBV	1079740	942663	16.73%	3.033%
6	6.469	743	745	752	rVB	93033	86265	1.53%	0.278%
7	6.810	799	803	806	rBV	1092757	871920	15.47%	2.805%
8	6.875	811	814	819	rBV	823436	655371	11.63%	2.108%
9	7.375	894	899	902	rBV	2552592	2660142	47.20%	8.558%
10	7.998	1001	1005	1008	rBV	1091662	920257	16.33%	2.961%
11	8.092	1017	1021	1025	rBV	1364623	1118704	19.85%	3.599%
12	8.722	1124	1128	1131	rBV	668045	551752	9.79%	1.775%
13	9.169	1199	1204	1208	rBV	3810589	4238542	75.21%	13.636%
14	9.286	1219	1224	1242	rVB2	172340	271081	4.81%	0.872%
15	9.563	1267	1271	1276	rBV	387338	357770	6.35%	1.151%
16	9.845	1315	1319	1328	rBV	1384719	1325344	23.52%	4.264%
17	10.633	1449	1453	1462	rBV	770992	853479	15.14%	2.746%
18	10.904	1493	1499	1502	rVB2	64142	61592	1.09%	0.198%
19	10.939	1502	1505	1510	rBV	112682	109461	1.94%	0.352%
20	11.198	1542	1549	1552	rBV6	39429	69396	1.23%	0.223%
21	11.333	1561	1572	1577	rBV	1431794	1485441	26.36%	4.779%
22	12.310	1734	1738	1741	rBV	351803	339996	6.03%	1.094%
23	12.563	1775	1781	1783	rBV3	91166	152614	2.71%	0.491%
24	12.592	1783	1786	1790	rVB3	62737	77104	1.37%	0.248%
25	12.668	1795	1799	1802	rBV2	155071	195809	3.47%	0.630%
26	12.715	1805	1807	1811	rVB2	57322	58213	1.03%	0.187%
27	12.927	1837	1843	1846	rBV	4307738	5635571	100.00%	18.130%
28	13.386	1918	1921	1926	rVB	63591	66613	1.18%	0.214%
29	13.439	1926	1930	1933	rBV	62110	62019	1.10%	0.200%
30	13.527	1942	1945	1949	rBV	407825	416280	7.39%	1.339%
31	13.562	1949	1951	1954	rVB	68822	57032	1.01%	0.183%
32	13.809	1990	1993	2000	rBV	772891	1028564	18.25%	3.309%
33	13.968	2014	2020	2026	rBV	1612043	1600761	28.40%	5.150%
34	14.145	2045	2050	2055	rBV4	37318	78414	1.39%	0.252%

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

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

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35	14.451	2095	2102	2105	rBV5	50495	112753	2.00%	0.363%
36	14.609	2126	2129	2134	rVB	57728	65022	1.15%	0.209%
37	15.415	2260	2266	2281	rBV	1159226	1689629	29.98%	5.436%
38	15.868	2338	2343	2348	rBV	40357	60129	1.07%	0.193%
39	15.927	2348	2353	2361	rBV2	76897	205358	3.64%	0.661%

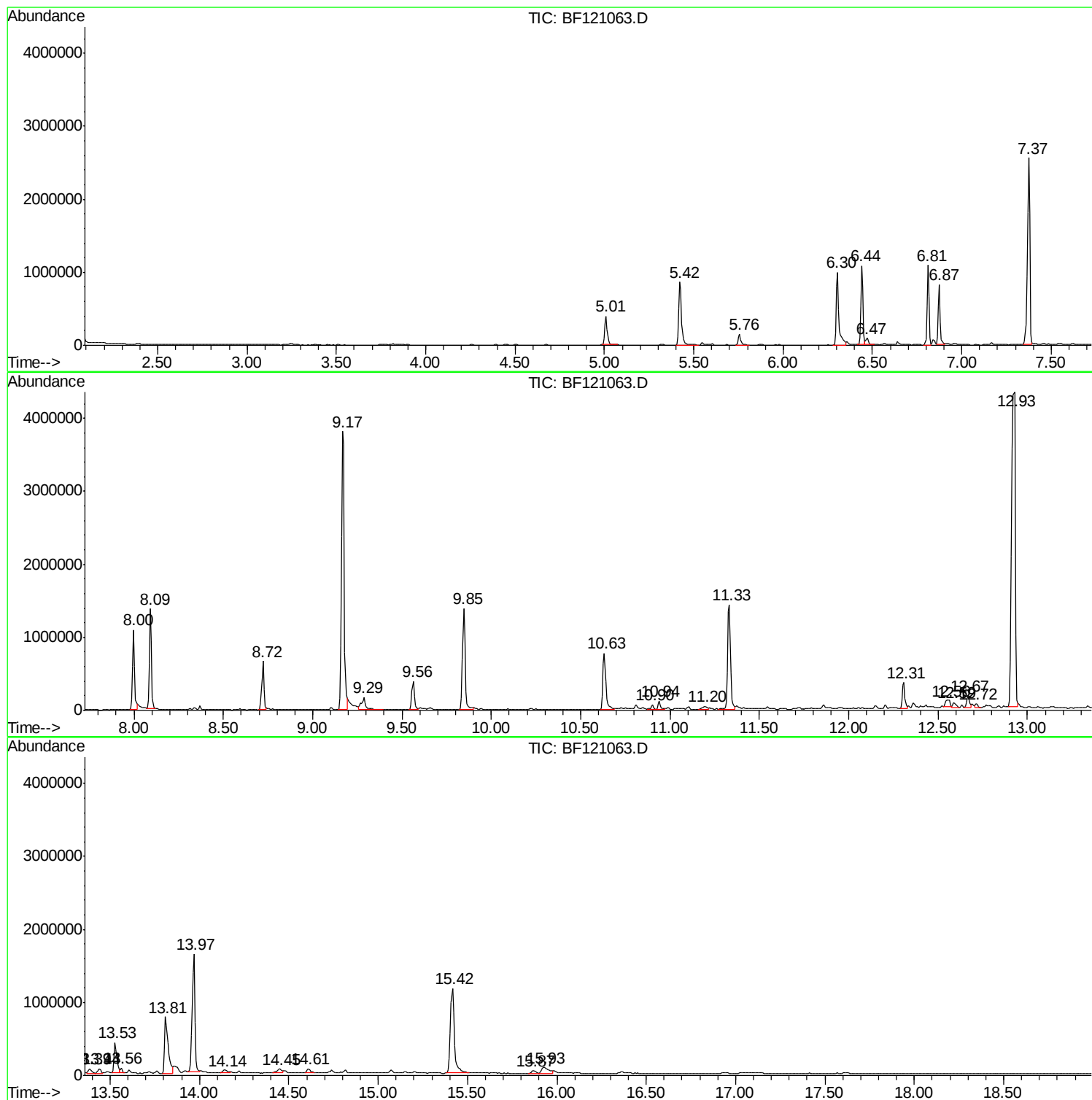
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071620.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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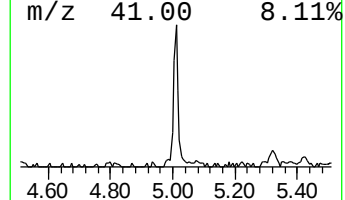
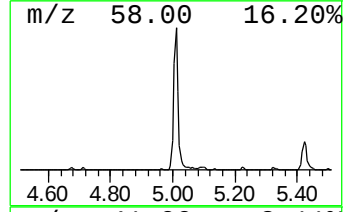
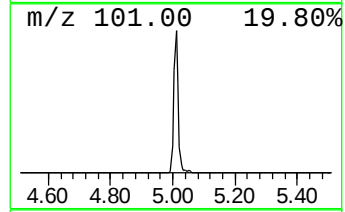
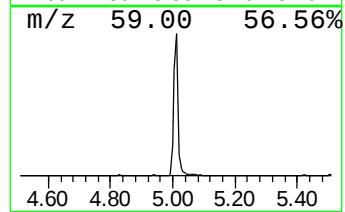
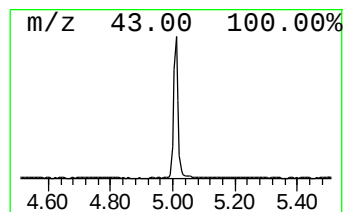
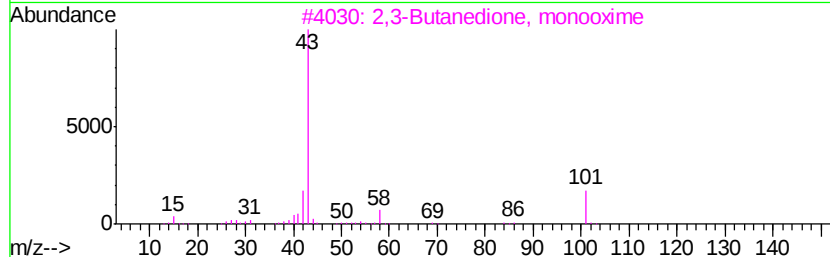
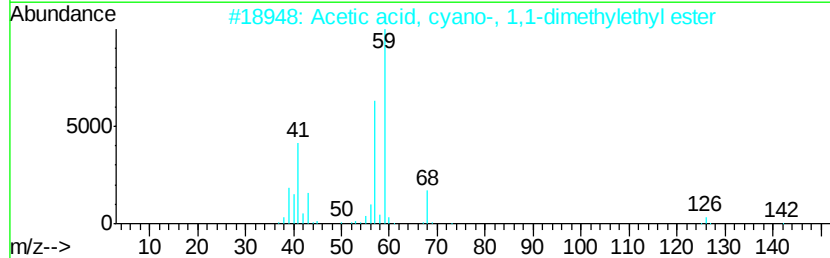
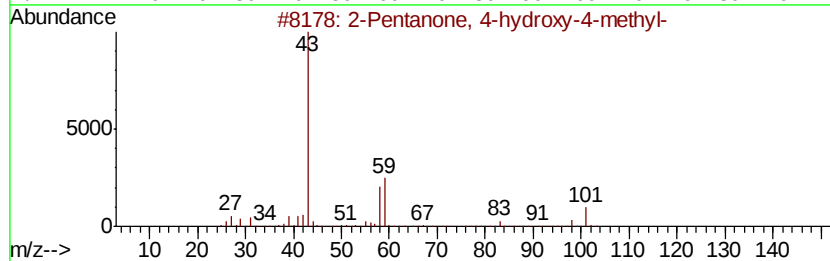
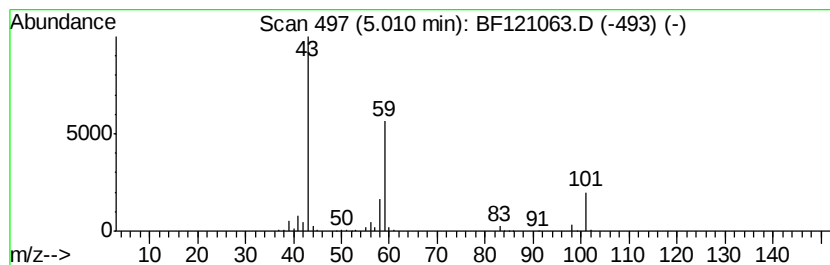
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.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.
5.01	8.77 ng	382257	1,4-Dichlorobenzene-d4	6.81

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	17
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
4		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	9
5		Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	9



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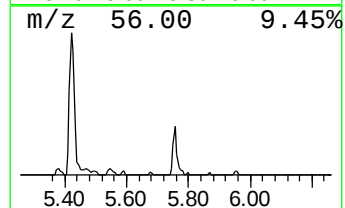
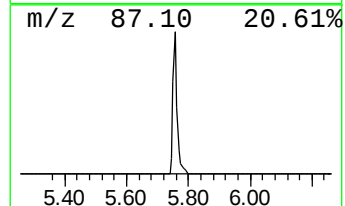
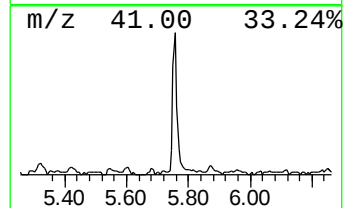
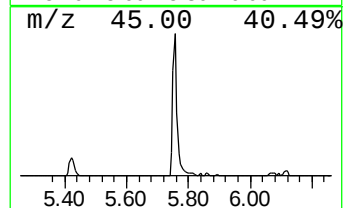
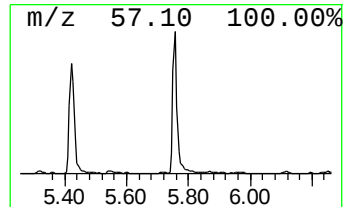
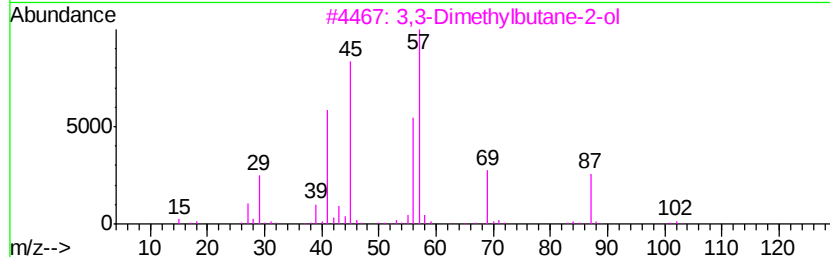
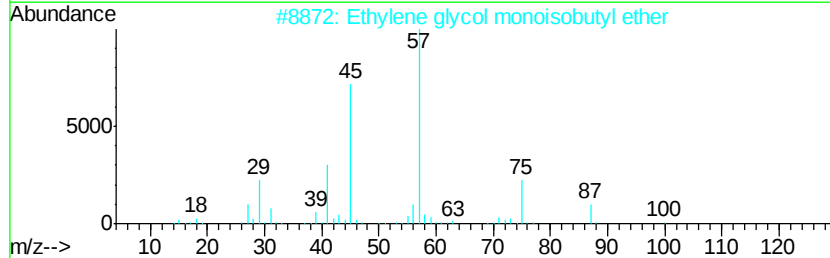
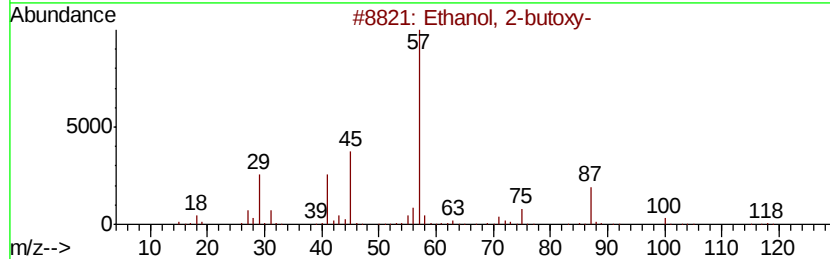
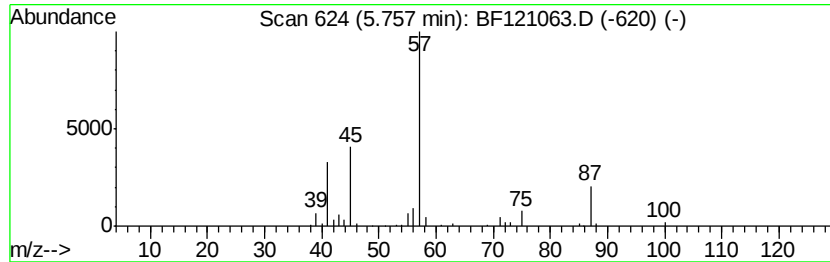
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Ethanol, 2-butoxy- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.76	3.21 ng	139980	1,4-Dichlorobenzene-d4	6.81

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Ethanol, 2-butoxy-	118	C6H14O2	000111-76-2	78
2		Ethylene glycol monoisobutyl ether	118	C6H14O2	004439-24-1	53
3		3,3-Dimethylbutane-2-ol	102	C6H14O	000464-07-3	42
4		n-Butyl ether	130	C8H18O	000142-96-1	25
5		1-Butene, 4-butoxy-	128	C8H16O	034061-76-2	23



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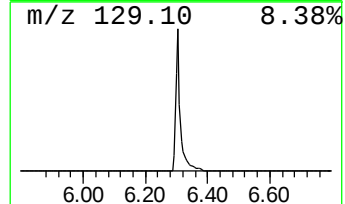
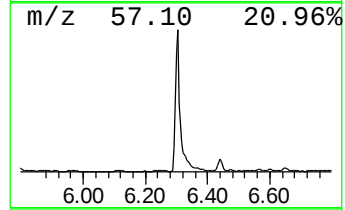
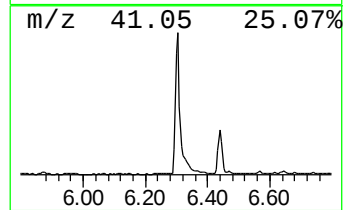
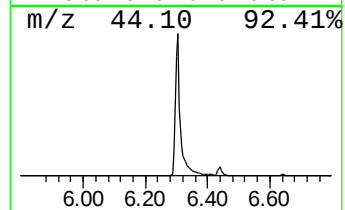
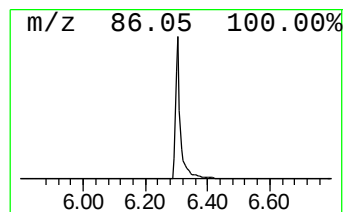
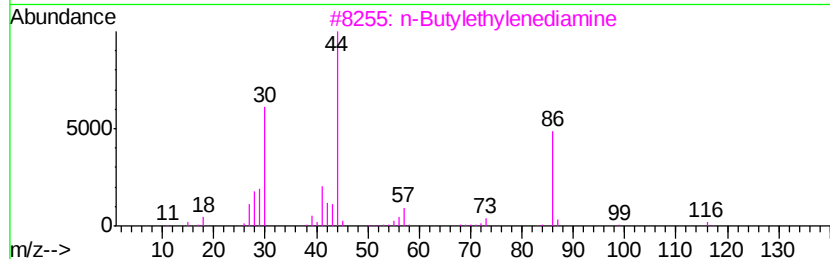
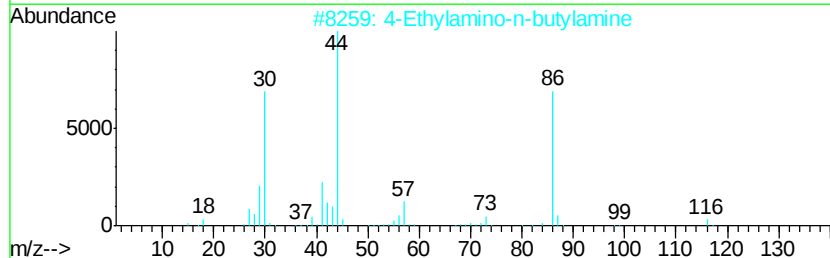
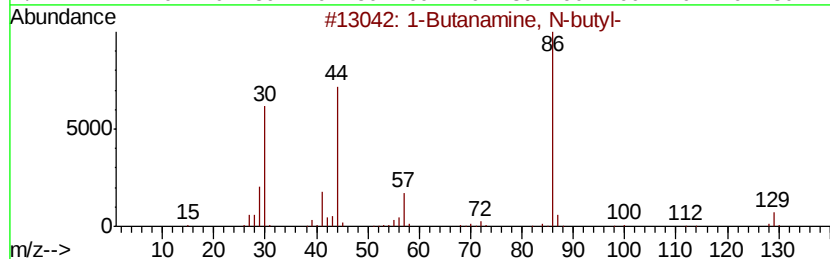
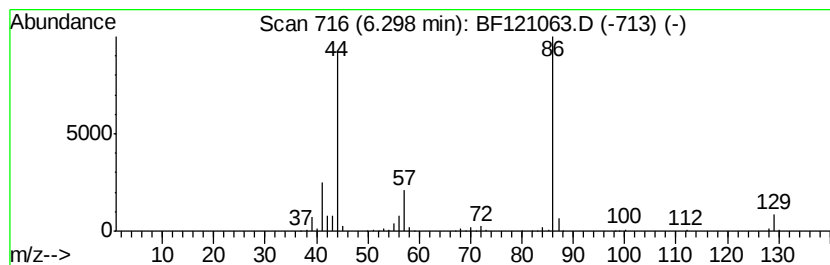
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 1-Butanamine, N-butyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.30	25.09 ng	1094030	1,4-Dichlorobenzene-d4	6.81

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Butanamine, N-butyl-	129	C8H19N	000111-92-2	91
2		4-Ethylamino-n-butylamine	116	C6H16N2	064429-16-9	86
3		n-Butylethylenediamine	116	C6H16N2	019522-69-1	64
4		1-Propanamine, 2-methyl-N-(2-met...	129	C8H19N	000110-96-3	49
5		1-Butanamine, N,N-diethyl-	129	C8H19N	004444-68-2	43



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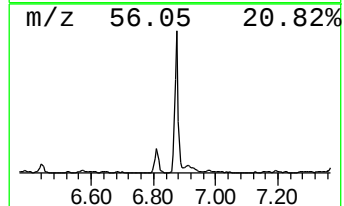
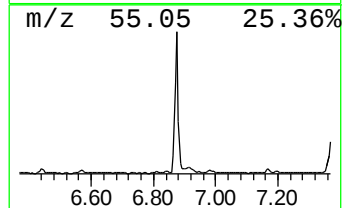
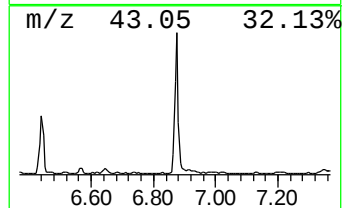
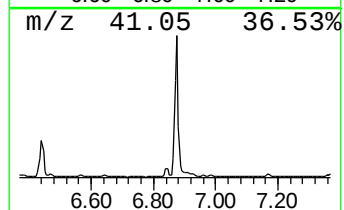
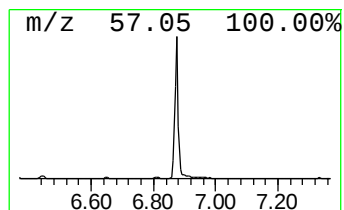
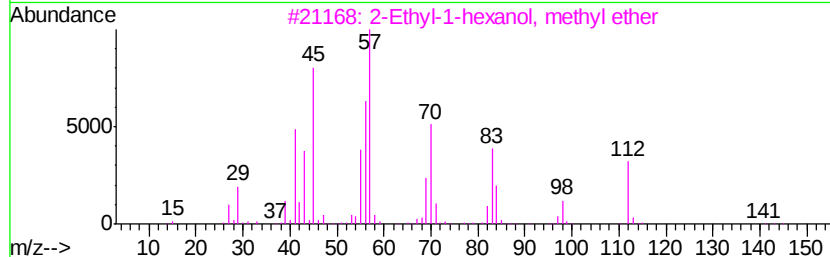
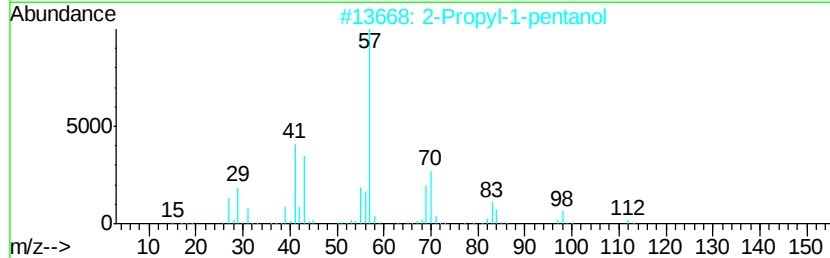
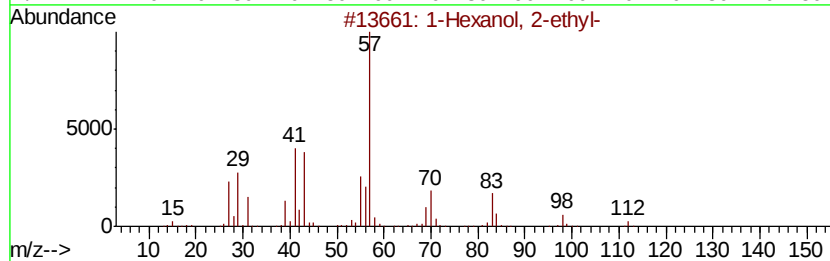
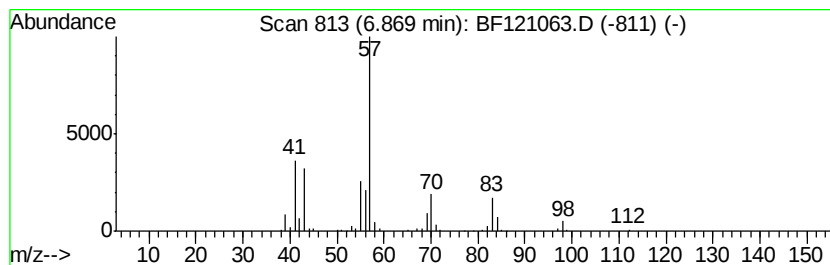
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 1-Hexanol, 2-ethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.87	15.03 ng	655371	1,4-Dichlorobenzene-d4	6.81

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	83
2		2-Propyl-1-pentanol	130	C8H18O	058175-57-8	53
3		2-Ethyl-1-hexanol, methyl ether	144	C9H20O	1000365-10-2	53
4		Heptane, 1,1'-oxybis-	214	C14H30O	000629-64-1	50
5		Oxalic acid, isobutyl octyl ester	258	C14H26O4	1000309-37-3	47



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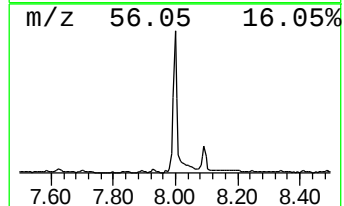
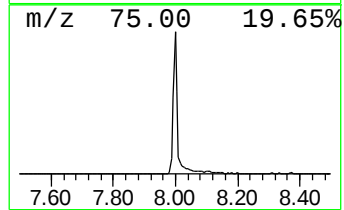
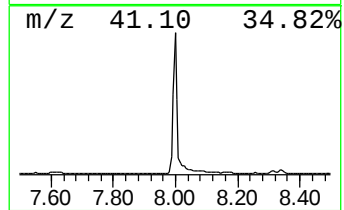
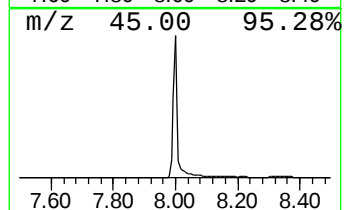
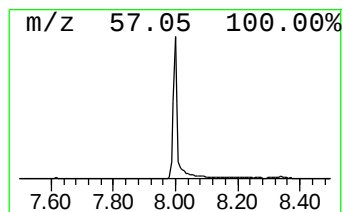
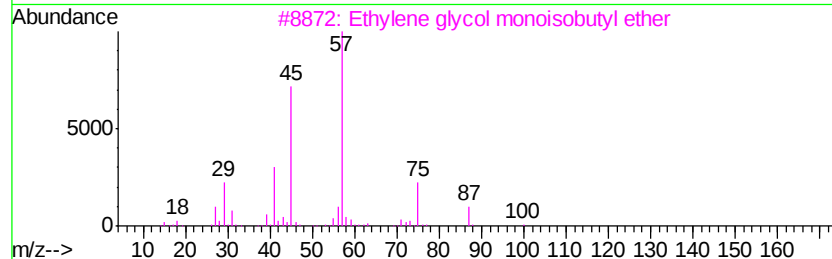
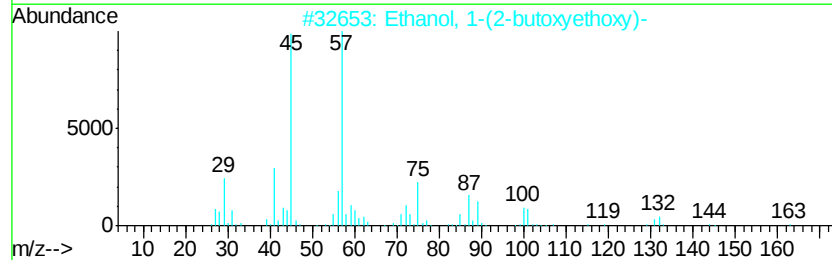
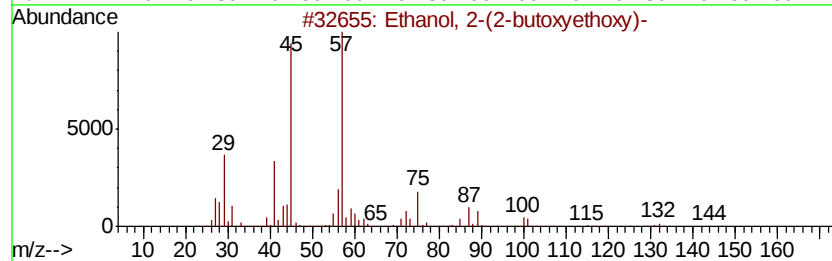
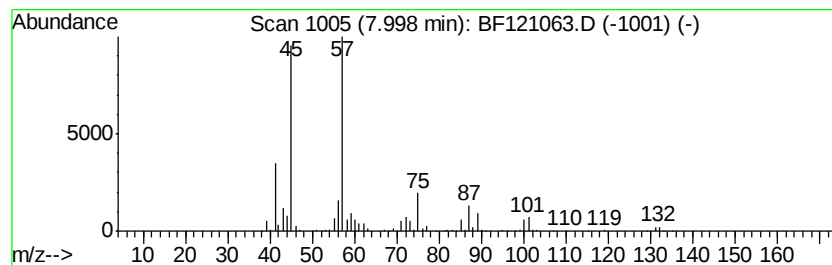
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.		
8.00	16.45 ng	920257	Naphthalene-d8	8.09		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
2		Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	83
3		Ethylene glycol monoisobutyl ether	118	C6H14O2	004439-24-1	64
4		3,3-Dimethylbutane-2-ol	102	C6H14O	000464-07-3	52
5		Methoxyacetic acid, 2-butyl ester	146	C7H14O3	070159-90-9	50



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072820\
Data File : BF121063.D
Acq On : 28 Jul 2020 11:48
Operator : JU/CG
Sample : L3378-16
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
DRUM-WATER

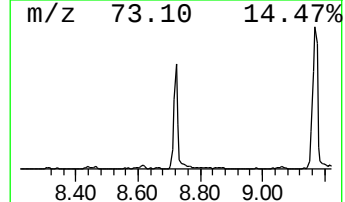
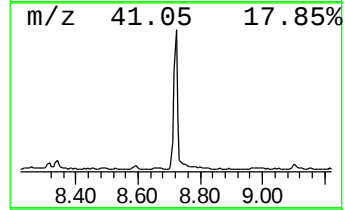
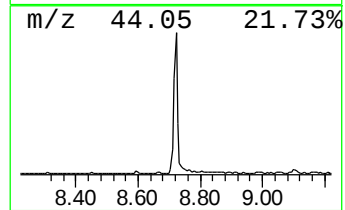
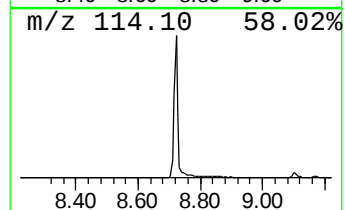
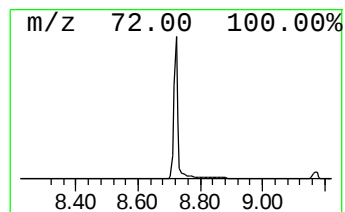
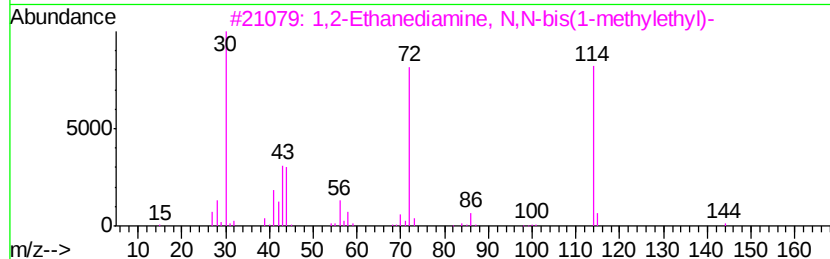
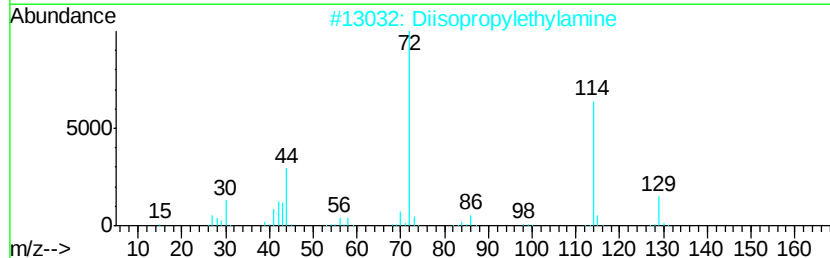
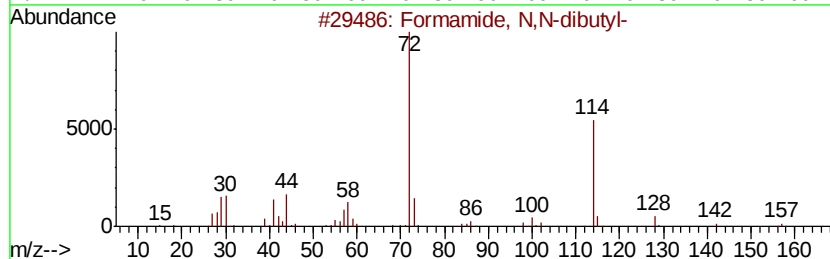
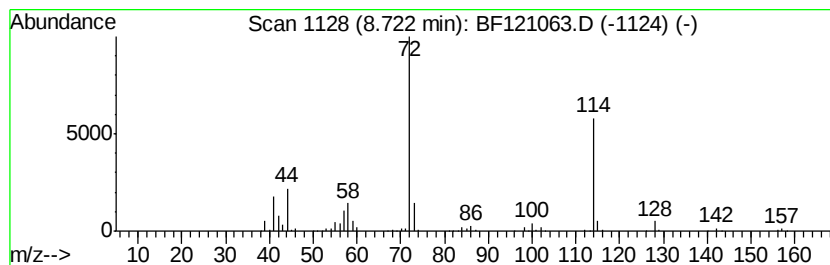
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Formamide, N,N-dibutyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.72	9.86 ng	551752	Napthalene-d8	8.09

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Formamide, N,N-dibutyl-	157	C9H19NO	000761-65-9	90
2		Diisopropylethylamine	129	C8H19N	007087-68-5	59
3		1,2-Ethanediamine, N,N-bis(1-met...	144	C8H20N2	000121-05-1	50
4		Isobutyl-propyl-amine	115	C7H17N	039190-66-4	38
5		Dimethyl(2-octyl)amine	157	C10H23N	007378-97-4	38



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072820\
Data File : BF121063.D
Acq On : 28 Jul 2020 11:48
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Sample : L3378-16
Misc :
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ClientSampleId :
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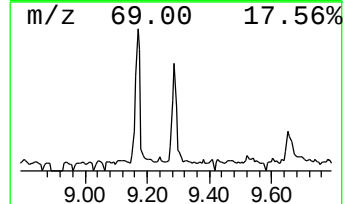
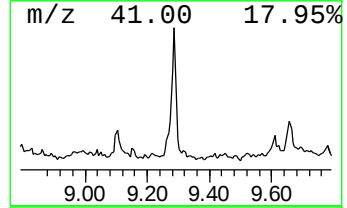
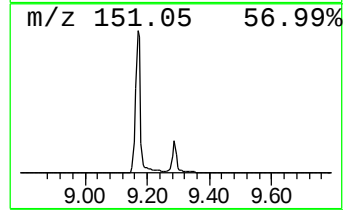
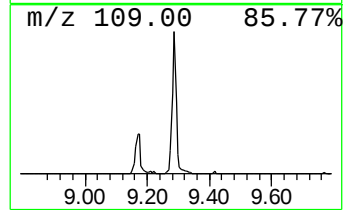
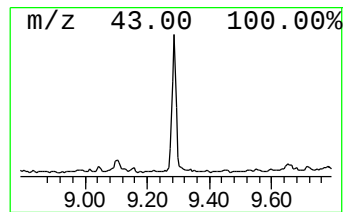
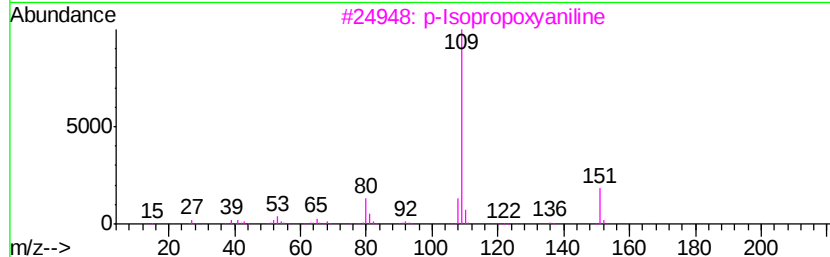
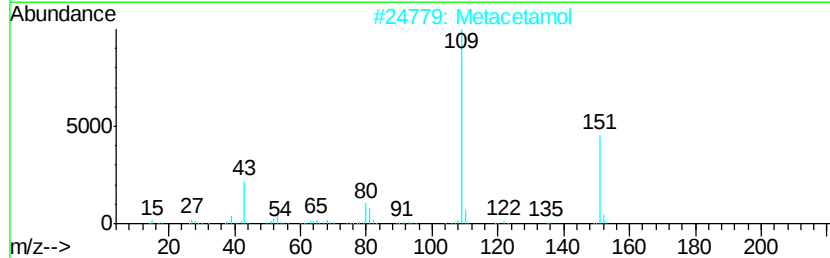
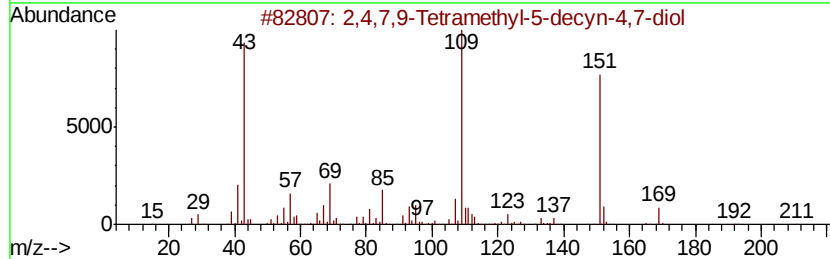
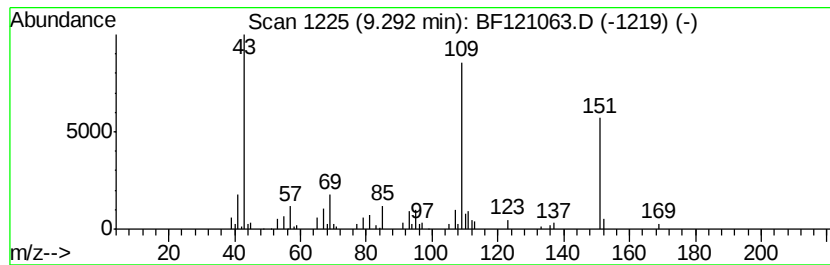
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 2,4,7,9-Tetramethyl-5-decyn... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.29	4.09 ng	271081	Acenaphthene-d10	9.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2,4,7,9-Tetramethyl-5-decyn-4,7-...	226	C14H26O2	000126-86-3	91
2		Metacetamol	151	C8H9NO2	000621-42-1	59
3		p-Isopropoxvaniline	151	C9H13NO	007664-66-6	59
4		1-(1,2,3-Trimethyl-cyclopent-2-e...	152	C10H16O	070987-81-4	46
5		Acetaminophen	151	C8H9NO2	000103-90-2	45



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072820\
Data File : BF121063.D
Acq On : 28 Jul 2020 11:48
Operator : JU/CG
Sample : L3378-16
Misc :
ALS Vial : 6 Sample Multiplier: 1

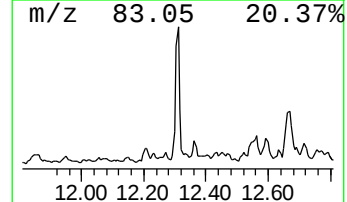
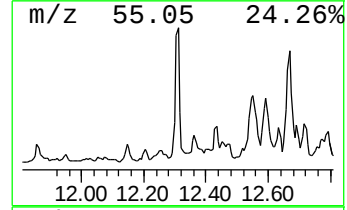
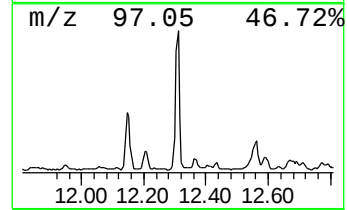
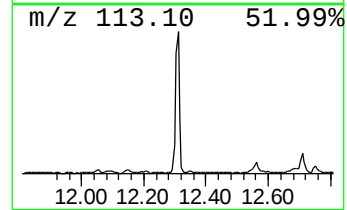
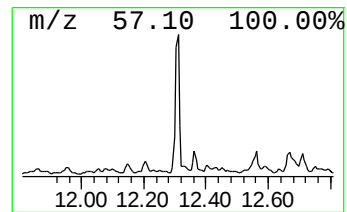
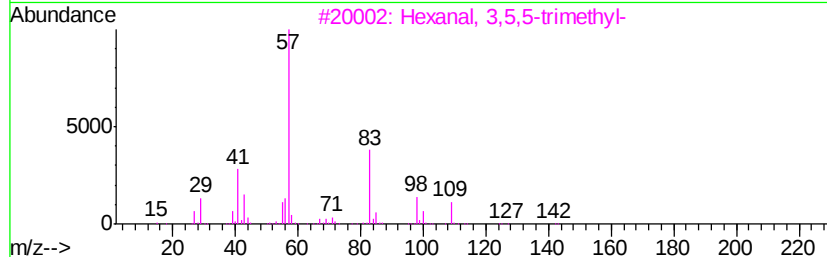
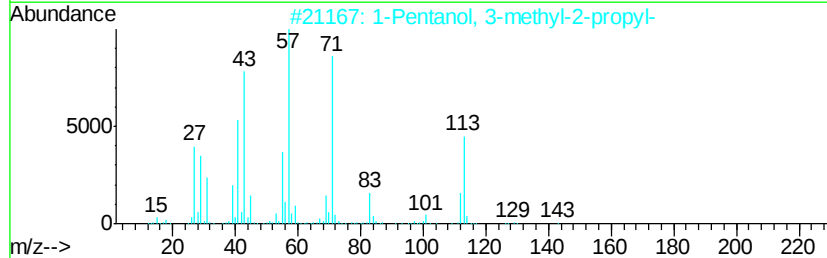
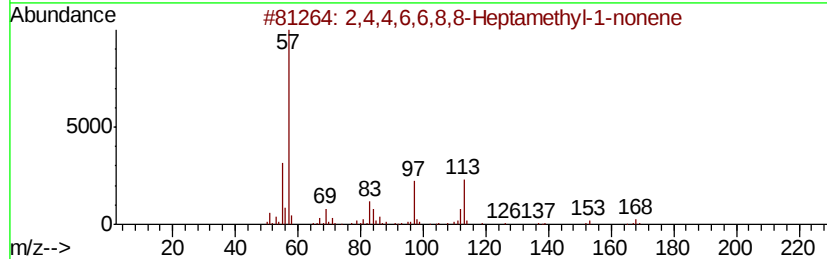
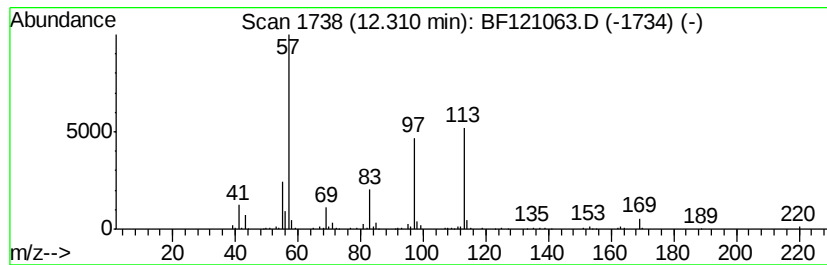
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071620.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 2,4,4,6,6,8,8-Heptamethyl-1... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.31	4.58 ng	339996	Phenanthrene-d10	11.33	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4,4,6,6,8,8-Heptamethyl-1-nonene	224	C16H32	015796-04-0	64
2	1-Pentanol, 3-methyl-2-propyl-	144	C9H20O	054004-40-9	38
3	Hexanal, 3,5,5-trimethyl-	142	C9H18O	005435-64-3	22
4	Vardenafil	488	C23H32N6O4S	224785-90-4	22
5	3-Methylpentyl ethylphosphonoflu...	196	C8H18FO2P	1000298-39-1	16



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072820\
Data File : BF121063.D
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Operator : JU/CG
Sample : L3378-16
Misc :
ALS Vial : 6 Sample Multiplier: 1

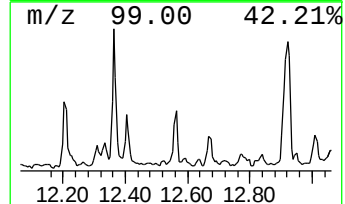
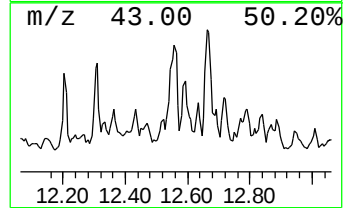
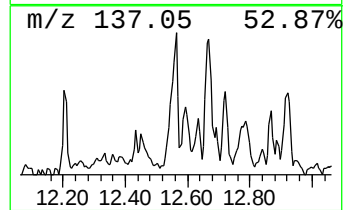
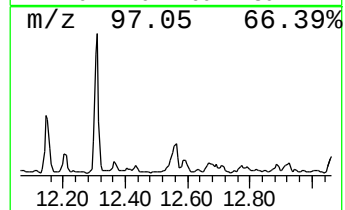
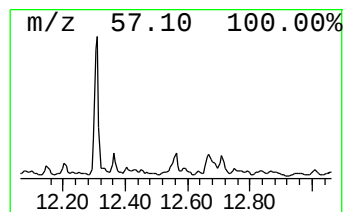
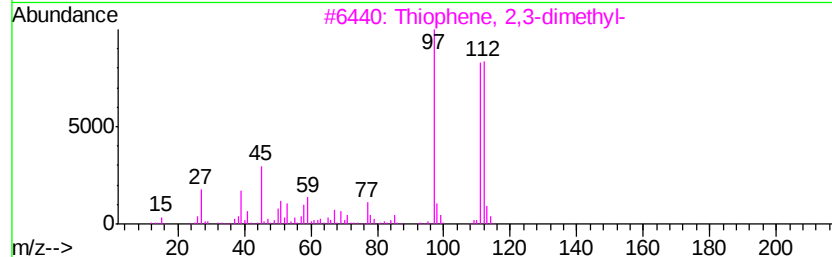
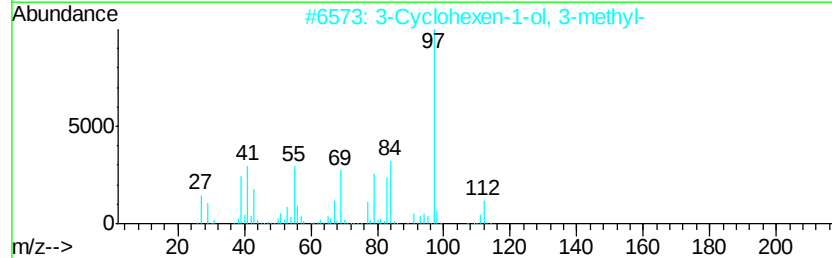
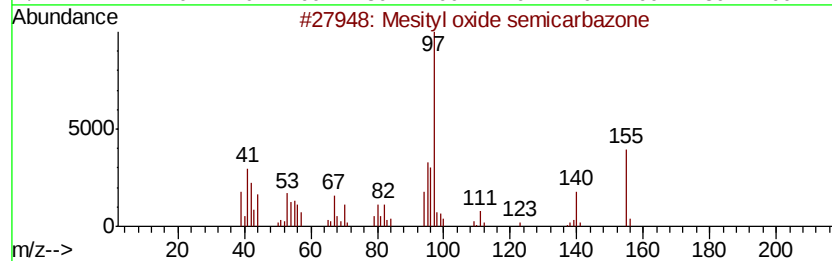
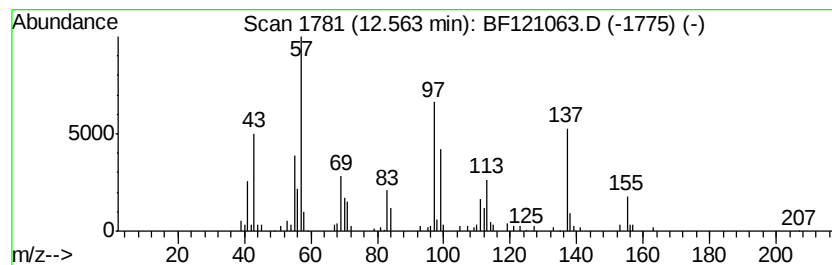
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071620.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 unknown12.56 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.56	2.05 ng	152614	Phenanthrene-d10	11.33
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Mesityl oxide semicarbazone	155 C7H13N3O	003780-62-9 18
2		3-Cyclohexen-1-ol, 3-methyl-	112 C7H12O	053783-91-8 14
3		Thiophene, 2,3-dimethyl-	112 C6H8S	000632-16-6 14
4		3-Pentenoic acid, 4-methyl-	114 C6H10O2	000504-85-8 11
5		2(5H)-Furanone, 5,5-dimethyl-	112 C6H8O2	020019-64-1 11



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072820\
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Sample : L3378-16
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
DRUM-WATER

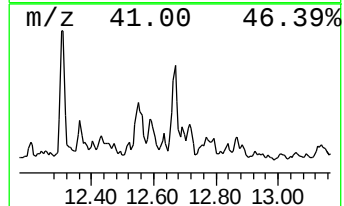
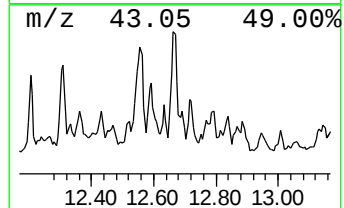
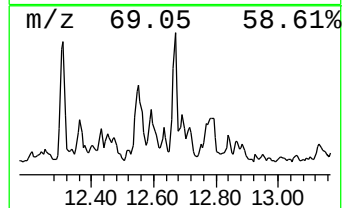
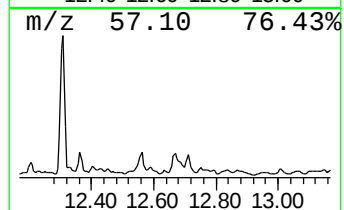
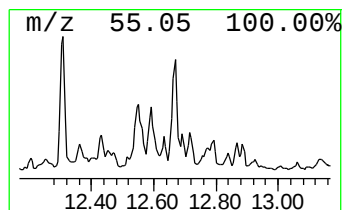
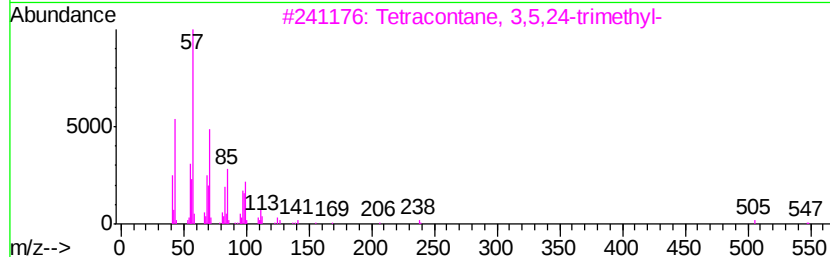
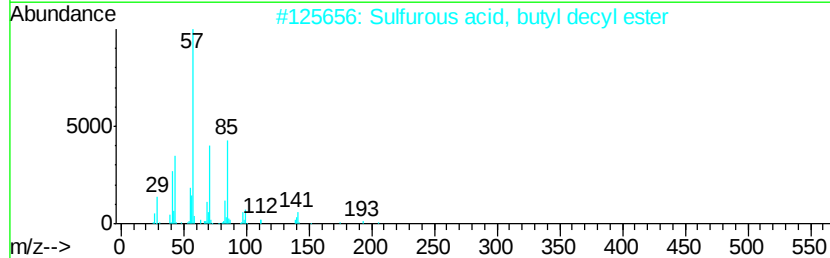
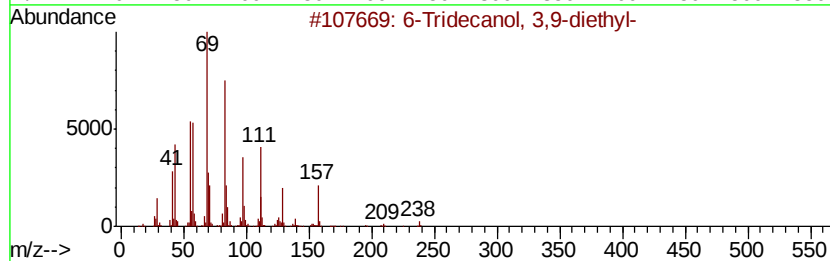
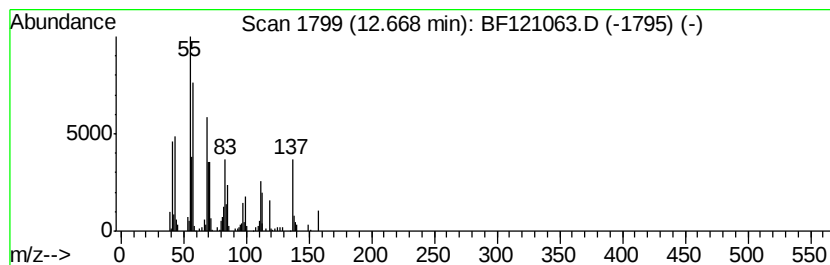
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 unknown12.67 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.67	2.45 ng	195809	Chrysene-d12	13.97

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	6-Tridecanol, 3,9-diethyl-			256	C17H36O	000123-24-0	35
2	Sulfurous acid, butyl decyl ester			278	C14H30O3S	1000309-17-7	27
3	Tetracontane, 3,5,24-trimethyl-			605	C43H88	055162-61-3	27
4	Tridecane, 1-bromo-			262	C13H27Br	000765-09-3	27
5	3-Ethyl-2-hexene			112	C8H16	000620-00-8	25



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072820\
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Misc :
ALS Vial : 6 Sample Multiplier: 1

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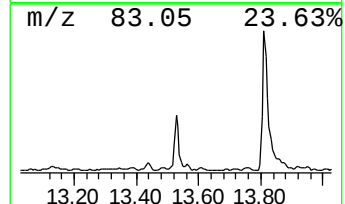
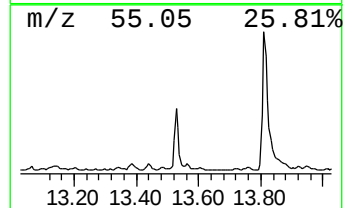
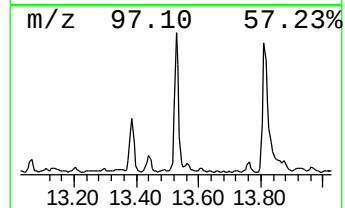
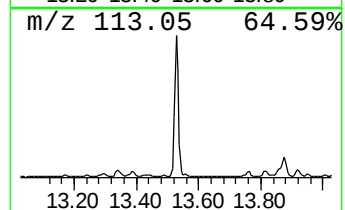
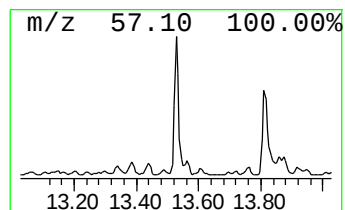
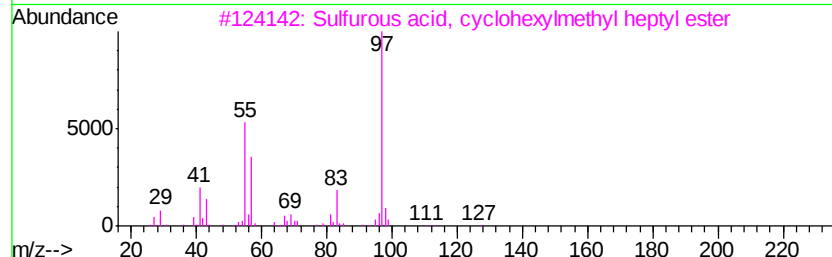
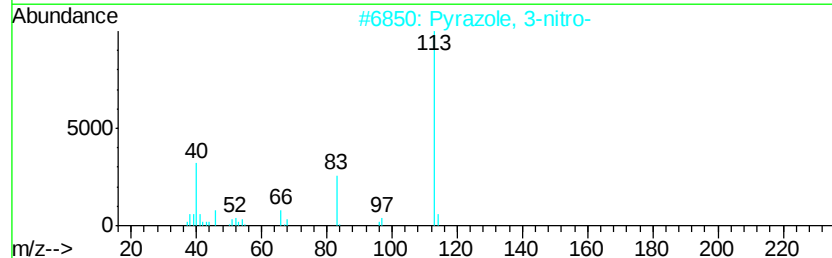
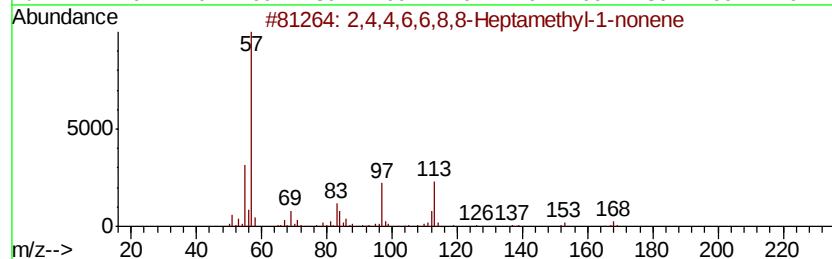
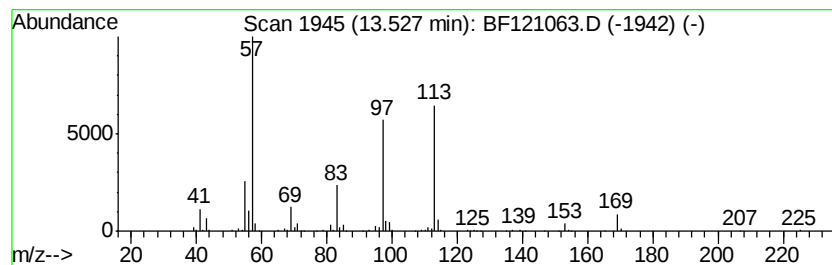
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 unknown13.53 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.53	5.20 ng	416280	Chrysene-d12	13.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2,4,4,6,6,8,8-Heptamethyl-1-nonene	224	C16H32	015796-04-0	50
2		Pyrazole, 3-nitro-	113	C3H3N3O2	026621-44-3	38
3		Sulfurous acid, cyclohexylmethyl...	276	C14H28O3S	1000309-21-7	38
4		2,4,4-Trimethyl-1-pentanol, pent...	276	C11H17F5O2	1000365-51-0	35
5		Cyclohexane, (bromomethyl)-	176	C7H13Br	002550-36-9	35



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072820\
Data File : BF121063.D
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Operator : JU/CG
Sample : L3378-16
Misc :
ALS Vial : 6 Sample Multiplier: 1

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ClientSampleId :
DRUM-WATER

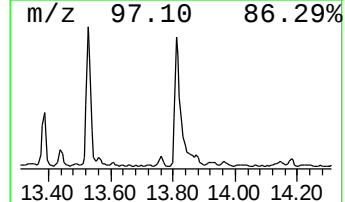
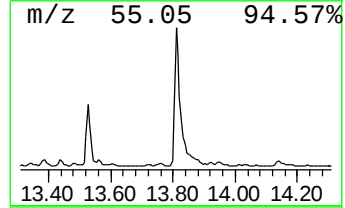
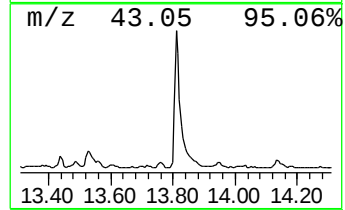
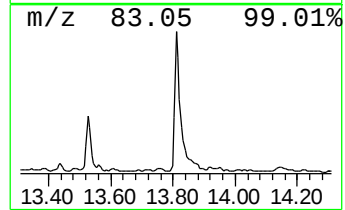
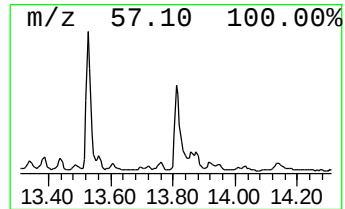
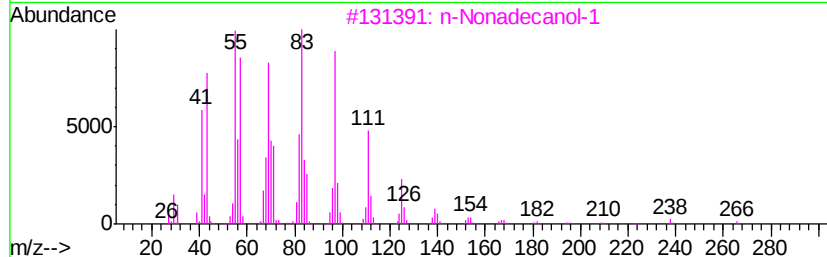
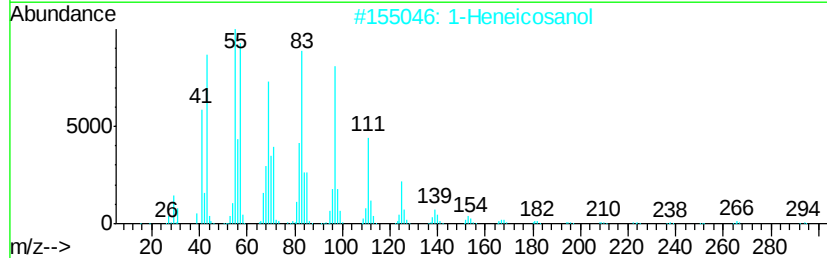
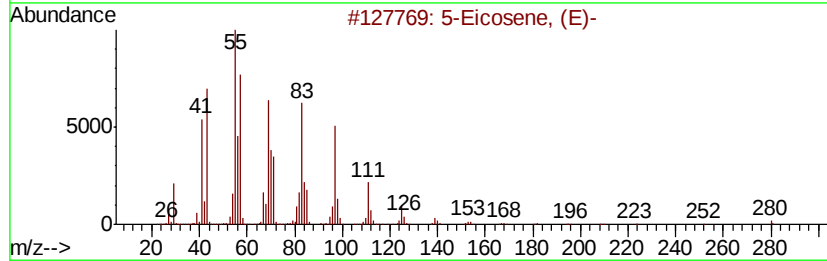
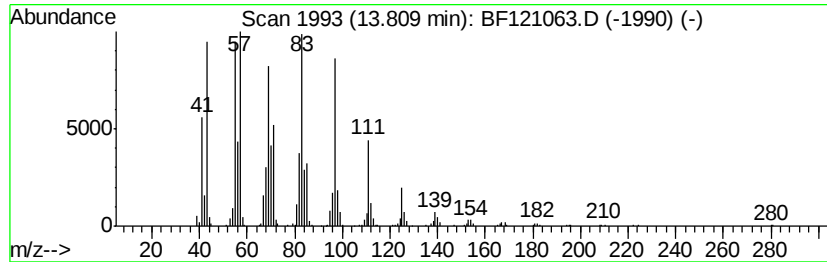
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 5-Eicosene, (E)- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.81	12.85 ng	1028560	Chrysene-d12	13.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Eicosene, (E)-	280	C20H40	074685-30-6	99
2		1-Heneicosanol	312	C21H44O	015594-90-8	95
3		n-Nonadecanol-1	284	C19H40O	001454-84-8	95
4		Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	94
5		Behenic alcohol	326	C22H46O	000661-19-8	94



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072820\
Data File : BF121063.D
Acq On : 28 Jul 2020 11:48
Operator : JU/CG
Sample : L3378-16
Misc :
ALS Vial : 6 Sample Multiplier: 1

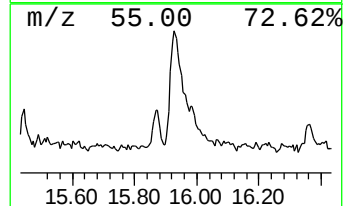
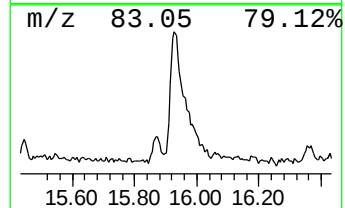
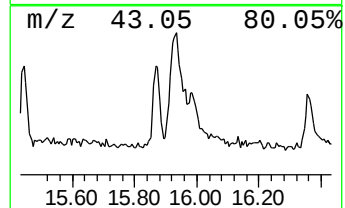
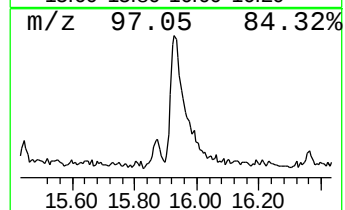
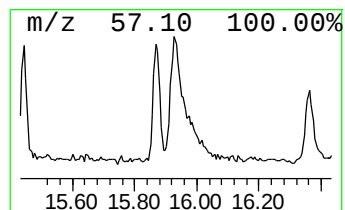
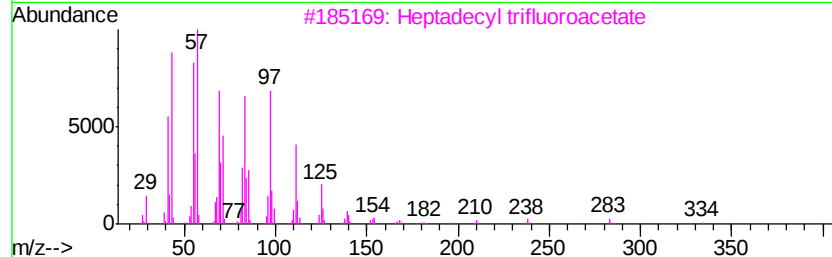
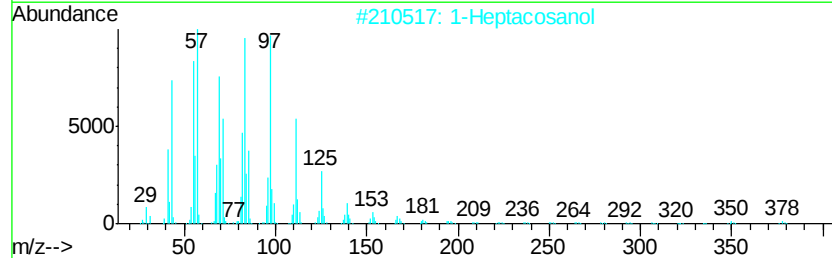
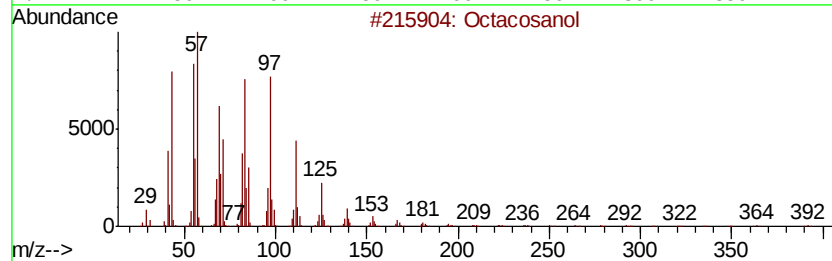
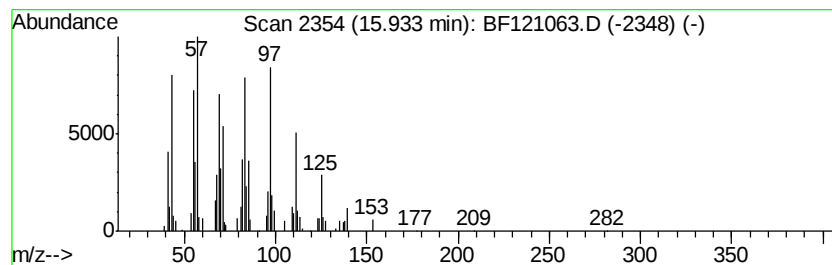
Instrument :
BNA_F
ClientSampleId :
DRUM-WATER

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071620.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Octacosanol Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.		
15.93	2.43 ng	205358	Perylene-d12	15.42		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octacosanol	410	C28H58O	000557-61-9	95
2		1-Heptacosanol	396	C27H56O	002004-39-9	95
3		Heptadecyl trifluoroacetate	352	C19H35F3O2	1000351-87-0	94
4		Heptafluorobutyric acid, n-octadec...	466	C22H37F7O2	000400-57-7	94
5		Heptafluorobutyric acid, hexadec...	438	C20H33F7O2	006385-15-5	94



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF072820\
 Data File : BF121063.D
 Acq On : 28 Jul 2020 11:48
 Operator : JU/CG
 Sample : L3378-16
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 DRUM-WATER

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF071620.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
2-Pentanone, 4-hy...	5.01	8.8	ng	382257	1	6.81	871920	20.0
Ethanol, 2-butoxy-	5.76	3.2	ng	139980	1	6.81	871920	20.0
1-Butanamine, N-b...	6.30	25.1	ng	1094030	1	6.81	871920	20.0
1-Hexanol, 2-ethyl-	6.87	15.0	ng	655371	1	6.81	871920	20.0
Ethanol, 2-(2-but...	8.00	16.4	ng	920257	2	8.09	1118700	20.0
Formamide, N,N-di...	8.72	9.9	ng	551752	2	8.09	1118700	20.0
2,4,7,9-Tetrameth...	9.29	4.1	ng	271081	3	9.85	1325340	20.0
2,4,4,6,6,8,8-Hep...	12.31	4.6	ng	339996	4	11.33	1485440	20.0
unknown12.56	12.56	2.0	ng	152614	4	11.33	1485440	20.0
unknown12.67	12.67	2.5	ng	195809	5	13.97	1600760	20.0
unknown13.53	13.53	5.2	ng	416280	5	13.97	1600760	20.0
5-Eicosene, (E)-	13.81	12.9	ng	1028560	5	13.97	1600760	20.0
Octacosanol	15.93	2.4	ng	205358	6	15.42	1689630	20.0