

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM082418\
 Data File : BM016594.D
 Acq On : 24 Aug 2018 18:39
 Operator : SJ/JU
 Sample : J4616-01
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 VB27184

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_M\METHODS\8270-BM082118.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	3.616	51	56	62	rBV	149060	210297	1.31%	0.310%
2	4.346	174	180	185	rBV2	111680	164069	1.02%	0.242%
3	5.387	351	357	362	rBV	109639	160420	1.00%	0.237%
4	5.551	381	385	399	rBV	553358	1006192	6.27%	1.485%
5	7.116	635	651	654	rBV2	124694	357691	2.23%	0.528%
6	7.192	660	664	675	rBV	276660	585430	3.65%	0.864%
7	7.534	716	722	727	rBV	1639767	2831901	17.65%	4.178%
8	7.881	775	781	791	rVB	180801	300557	1.87%	0.443%
9	8.004	796	802	810	rBV	457287	806271	5.03%	1.190%
10	8.116	810	821	827	rVB	4234487	6603336	41.17%	9.742%
11	8.322	851	856	870	rVV	1948587	3253788	20.28%	4.801%
12	8.486	873	884	892	rVB4	95787	251340	1.57%	0.371%
13	9.081	978	985	993	rBV7	86334	174926	1.09%	0.258%
14	9.192	998	1004	1013	rBV	1152046	1998882	12.46%	2.949%
15	9.481	1038	1053	1060	rVB6	167989	457121	2.85%	0.674%
16	9.822	1107	1111	1123	rVB4	131652	251791	1.57%	0.371%
17	10.810	1271	1279	1285	rBV	564574	1044016	6.51%	1.540%
18	11.957	1470	1474	1484	rVB2	114138	213744	1.33%	0.315%
19	13.263	1689	1696	1710	rBV	5555802	7694993	47.97%	11.353%
20	14.110	1835	1840	1853	rBV	674321	1127122	7.03%	1.663%
21	14.645	1926	1931	1941	rVB2	1020224	1612723	10.05%	2.379%
22	16.139	2180	2185	2195	rBV	6077796	8376065	52.22%	12.358%
23	17.392	2393	2398	2405	rBV	1767285	2487376	15.51%	3.670%
24	18.251	2540	2544	2554	rBV	980371	1410331	8.79%	2.081%
25	19.515	2755	2759	2769	rBV2	547468	792196	4.94%	1.169%
26	19.980	2833	2838	2844	rBV	14584386	16040727	100.00%	23.666%
27	21.215	3045	3048	3054	rVB2	542264	663256	4.13%	0.979%
28	21.533	3098	3102	3107	rBV	2854967	3432288	21.40%	5.064%
29	23.815	3484	3490	3500	rVB	1725454	3469878	21.63%	5.119%

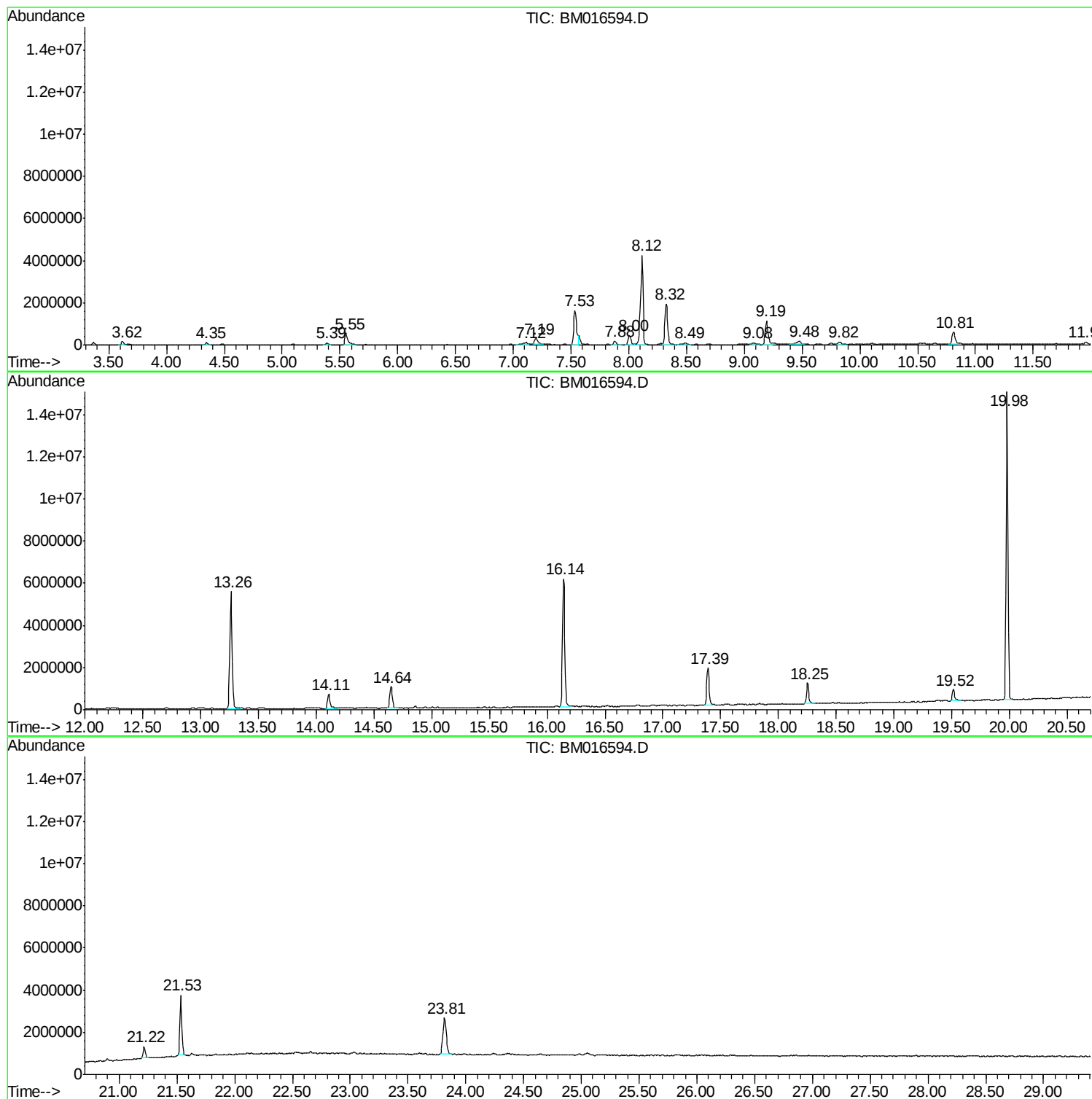
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM082118.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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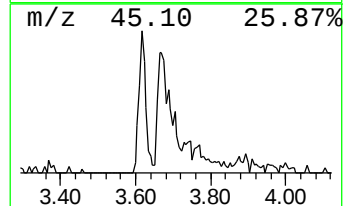
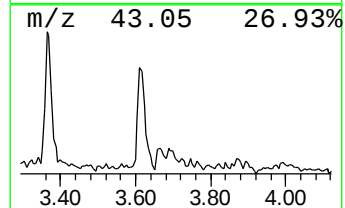
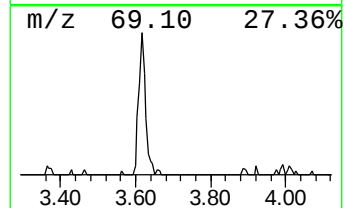
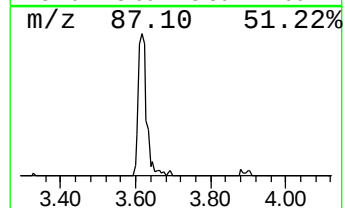
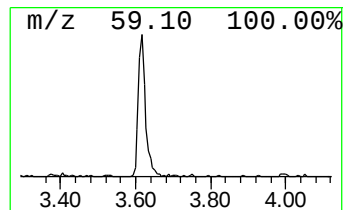
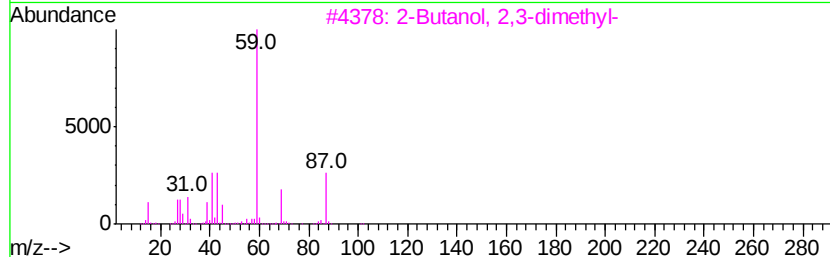
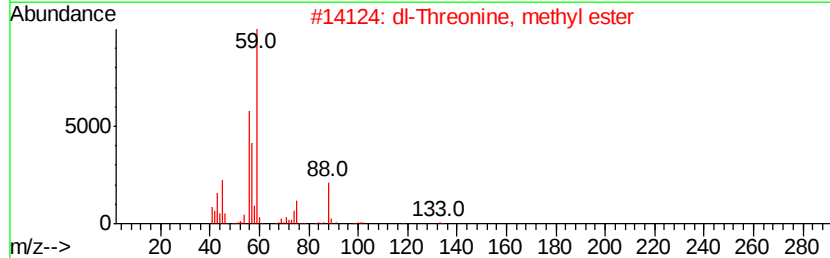
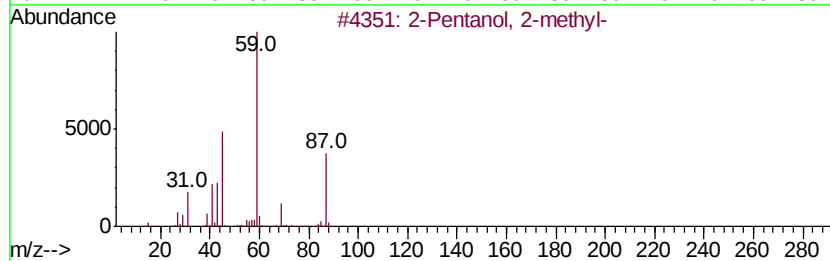
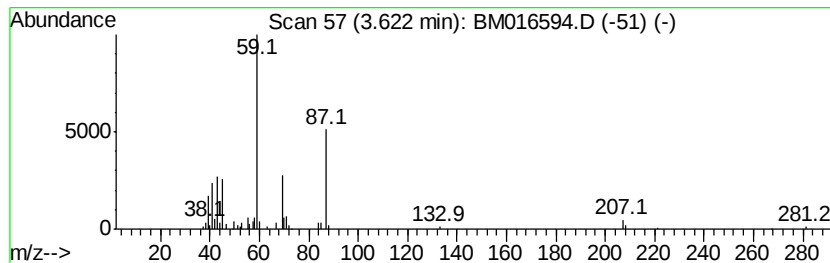
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM082118.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanol, 2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.62	5.22 ng	210297	1,4-Dichlorobenzene-d4	8.01

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanol, 2-methyl-	102	C6H14O	000590-36-3	64
2			dl-Threonine, methyl ester	133	C5H11NO3	1000130-33-2	53
3			2-Butanol, 2,3-dimethyl-	102	C6H14O	000594-60-5	50
4			4-Pentene-2-ol, 2-methyl	100	C6H12O	000624-97-5	47
5			Propane, 2-ethoxy-2-methyl-	102	C6H14O	000637-92-3	38



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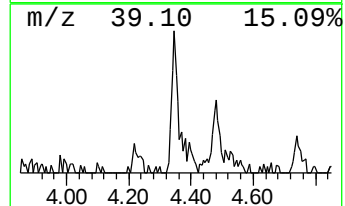
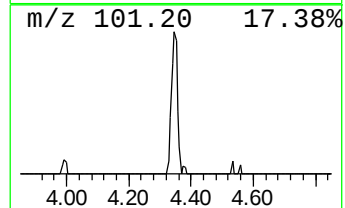
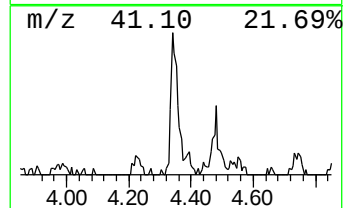
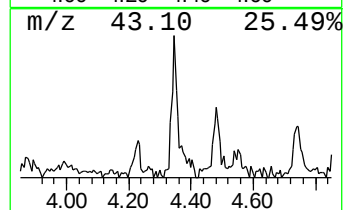
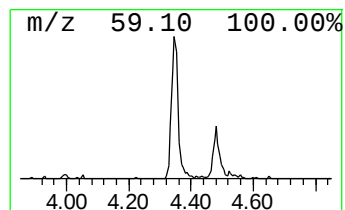
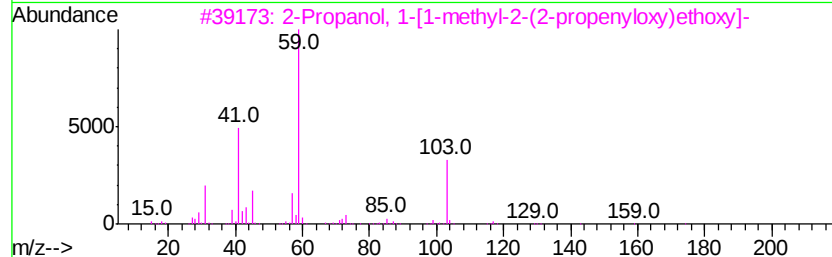
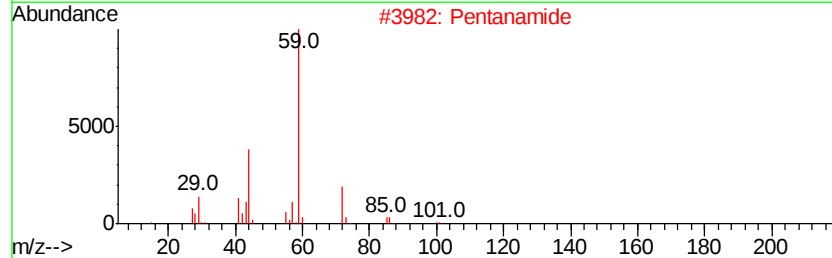
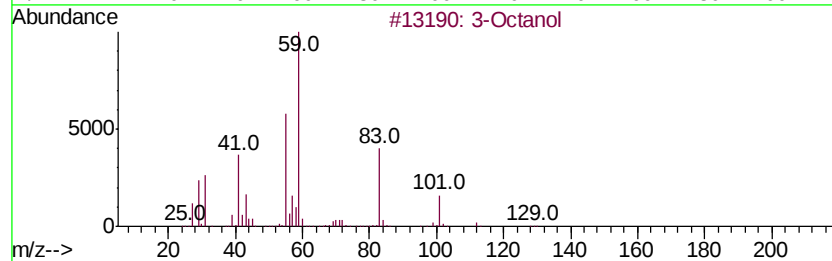
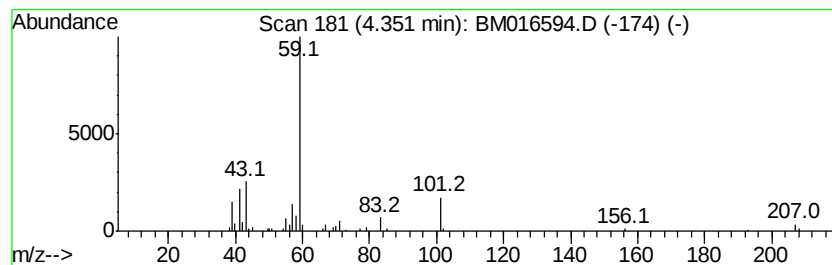
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Peak Number 2 3-Octanol Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.35	4.07 ng	164069	1,4-Dichlorobenzene-d4	8.01

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Octanol	130	C8H18O	000589-98-0	64
2			Pentanamide	101	C5H11NO	000626-97-1	59
3			2-Propanol, 1-[1-methyl-2-(2-pro...	174	C9H18O3	055956-25-7	56
4			2-Pentanol, 2,4-dimethyl-	116	C7H16O	000625-06-9	56
5			3-Octanol	130	C8H18O	020296-29-1	56



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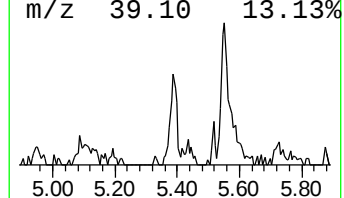
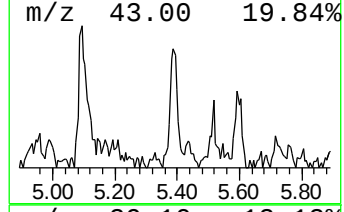
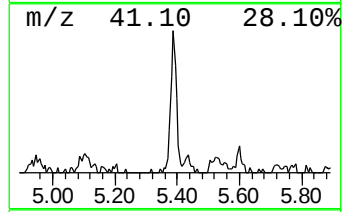
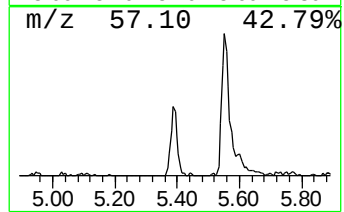
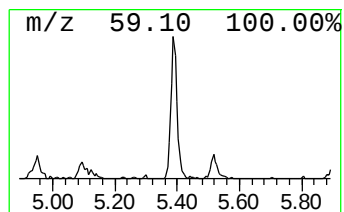
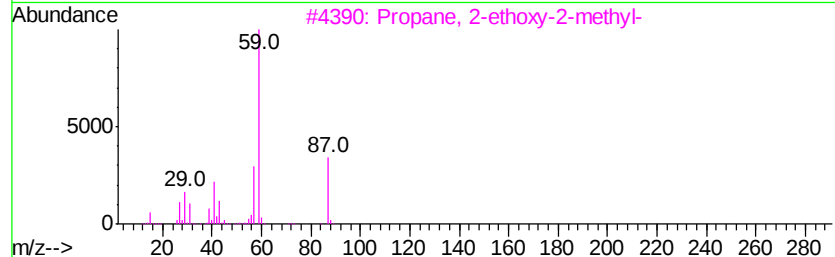
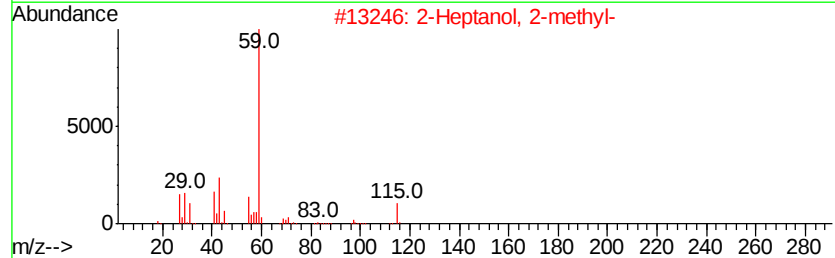
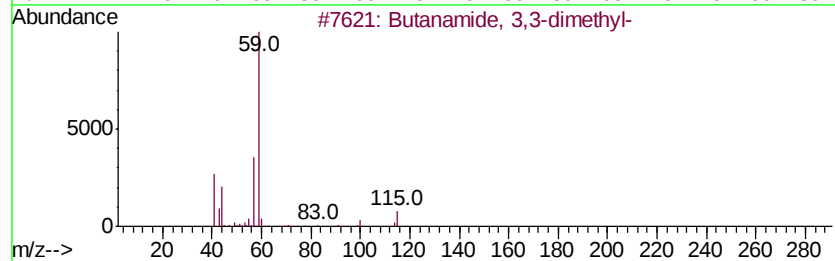
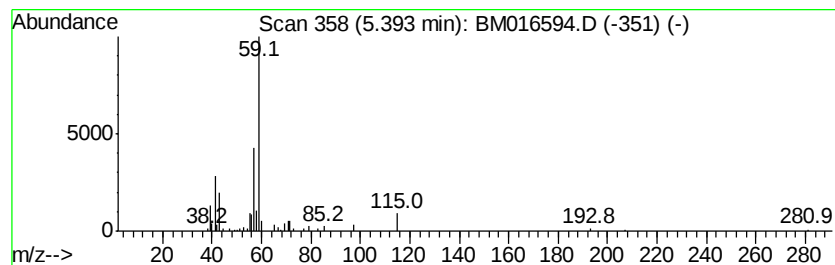
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Peak Number 3 Butanamide, 3,3-dimethyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.39	3.98 ng	160420	1,4-Dichlorobenzene-d4	8.01

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butanamide, 3,3-dimethyl-	115	C6H13NO	000926-04-5	72
2		2-Heptanol, 2-methyl-	130	C8H18O	000625-25-2	39
3		Propane, 2-ethoxy-2-methyl-	102	C6H14O	000637-92-3	38
4		t-Butyl ethyl ether	102	C6H14O	1000118-18-4	38
5		1-Butanol, 3-methoxy-	104	C5H12O2	002517-43-3	38



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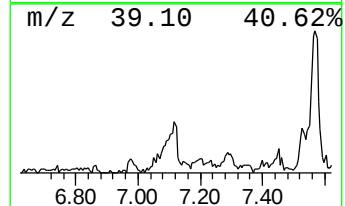
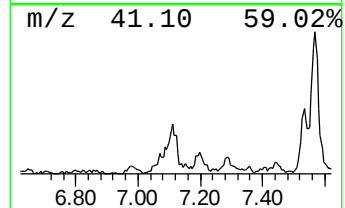
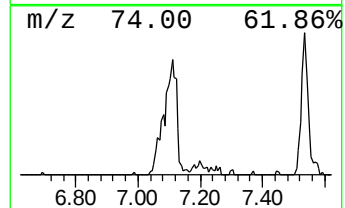
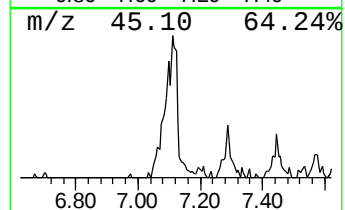
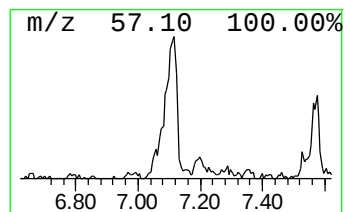
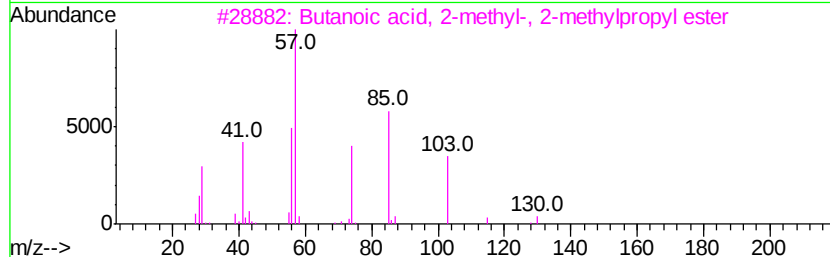
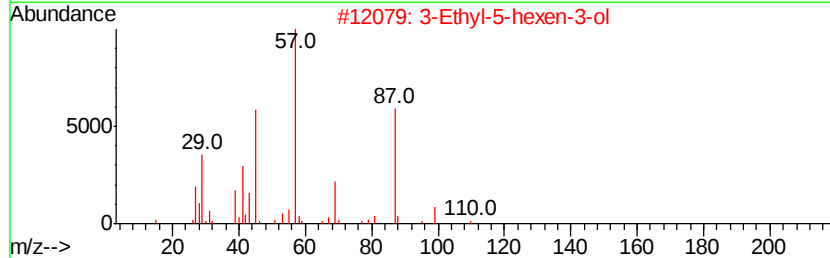
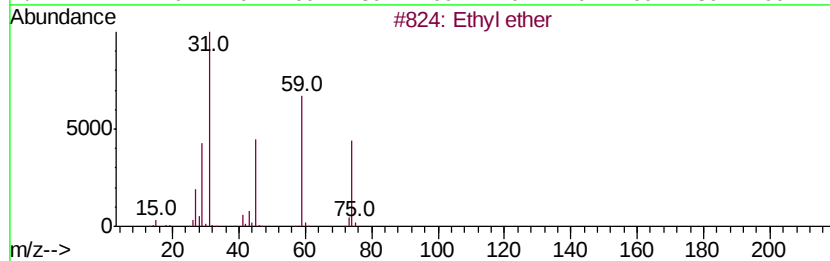
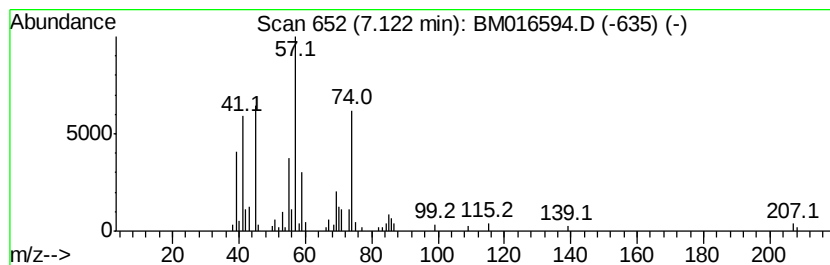
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 4 unknown7.12 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.12	8.87 ng	357691	1,4-Dichlorobenzene-d4	8.01

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Ethyl ether	74	C4H10O	000060-29-7	37
2		3-Ethyl-5-hexen-3-ol	128	C8H16O	001907-46-6	16
3		Butanoic acid, 2-methyl-, 2-meth...	158	C9H18O2	002445-67-2	12
4		Methoxyacetic acid, 2-tridecyl e...	272	C16H32O3	1000282-04-5	12
5		2H-Pyran, 2-(1,1-dimethylethoxy)...	158	C9H18O2	001927-69-1	10



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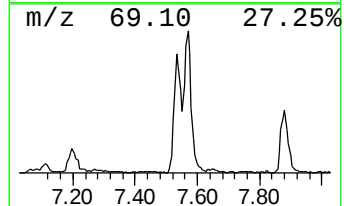
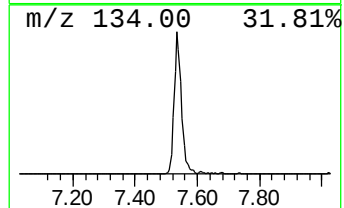
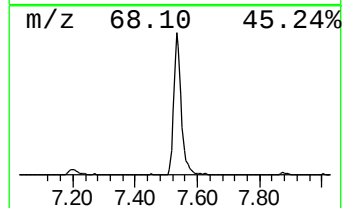
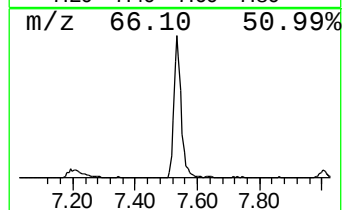
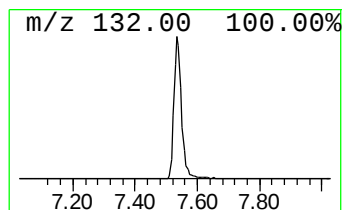
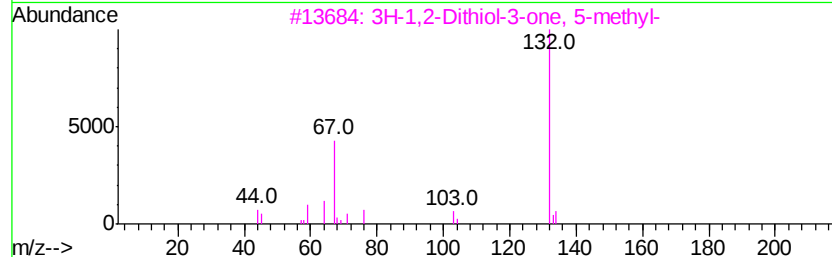
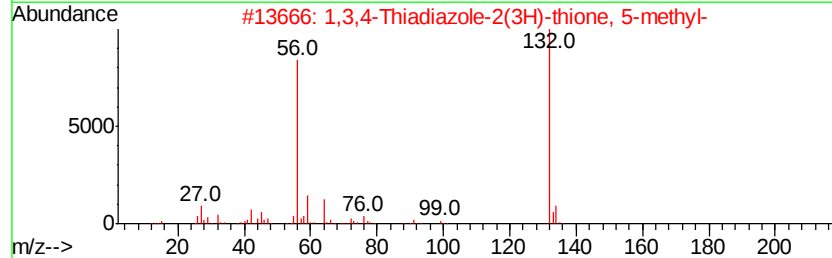
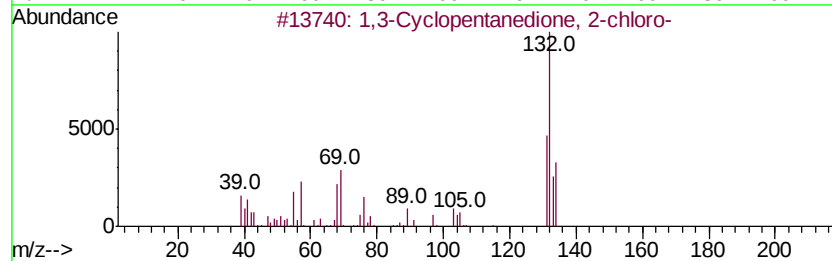
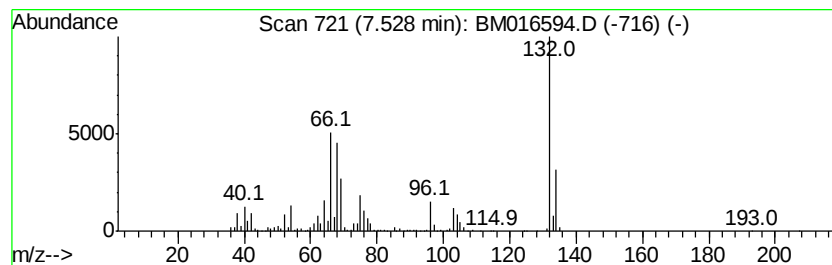
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 5 unknown7.53 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.53	70.25 ng	2831900	1,4-Dichlorobenzene-d4	8.01

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	10
2		1,3,4-Thiadiazole-2(3H)-thione, ...	132	C3H4N2S2	029490-19-5	10
3		3H-1,2-Dithiol-3-one, 5-methyl-	132	C4H4OS2	003620-08-4	10
4		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	10
5		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9



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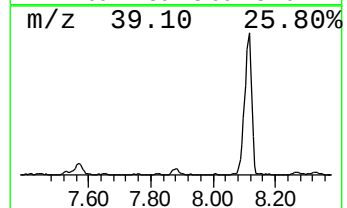
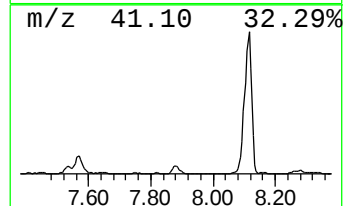
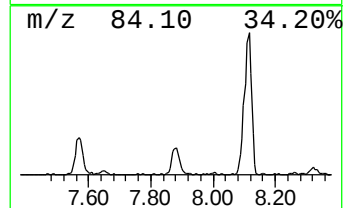
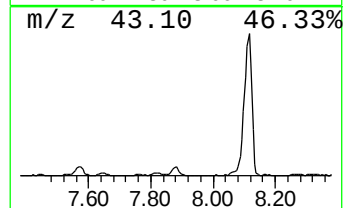
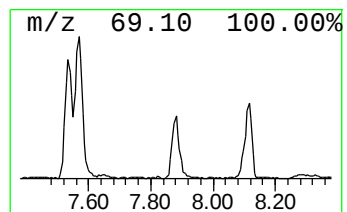
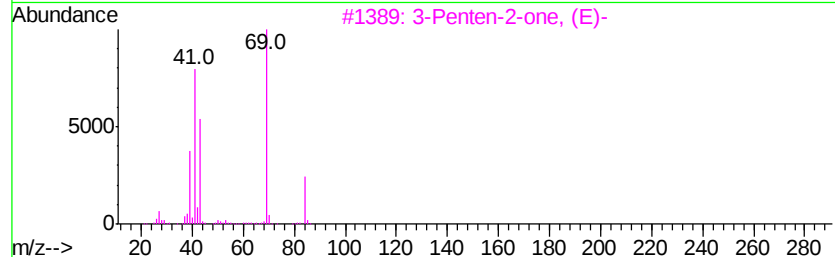
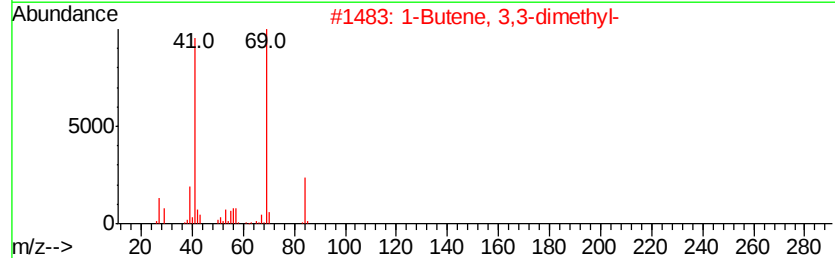
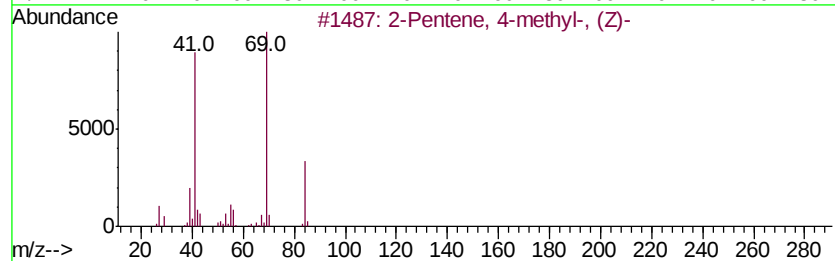
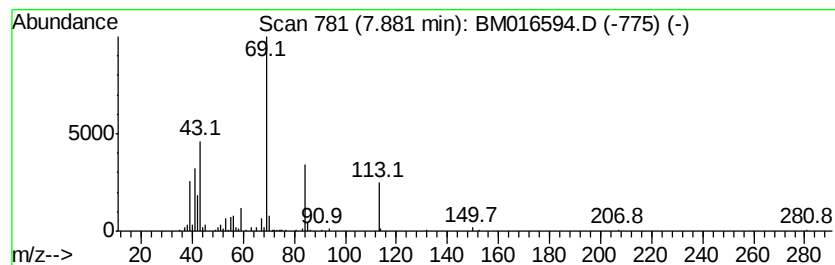
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 2-Pentene, 4-methyl-, (Z)- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.88	7.46 ng	300557	1,4-Dichlorobenzene-d4	8.01

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentene, 4-methyl-, (Z)-	84	C6H12	000691-38-3	58
2			1-Butene, 3,3-dimethyl-	84	C6H12	000558-37-2	53
3			3-Penten-2-one, (E)-	84	C5H8O	003102-33-8	53
4			3-Penten-2-one	84	C5H8O	000625-33-2	53
5			Furan, 2,3-dihydro-3-methyl-	84	C5H8O	001708-27-6	50



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM082418\
Data File : BM016594.D
Acq On : 24 Aug 2018 18:39
Operator : SJ/JU
Sample : J4616-01
Misc :
ALS Vial : 17 Sample Multiplier: 1

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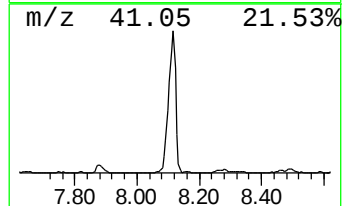
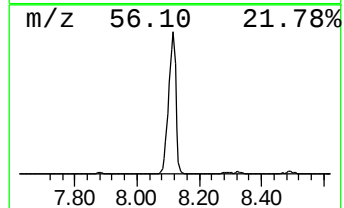
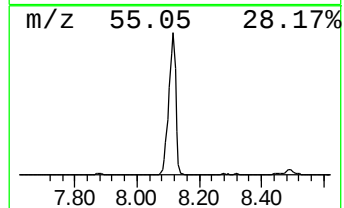
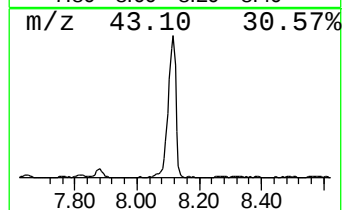
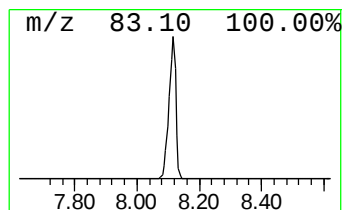
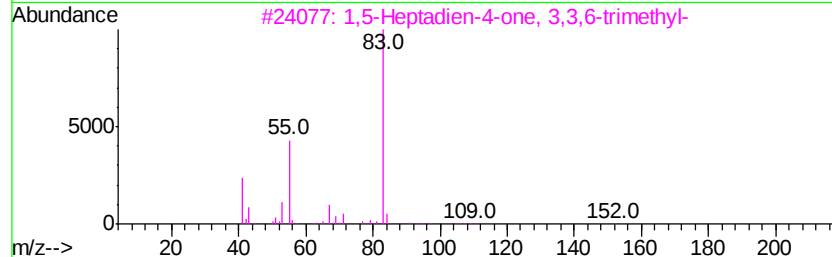
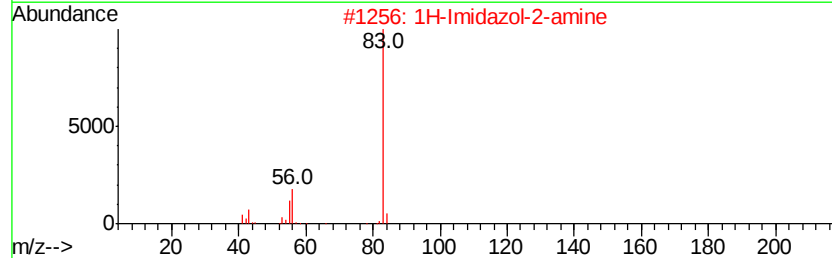
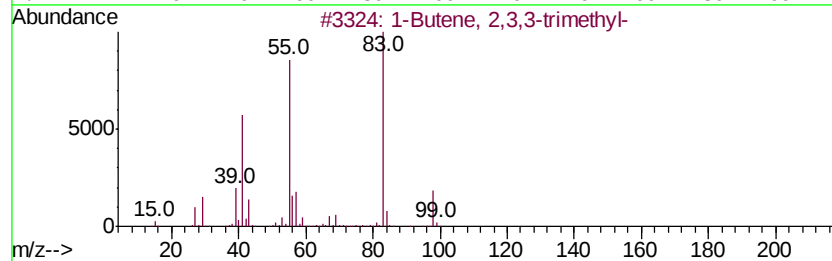
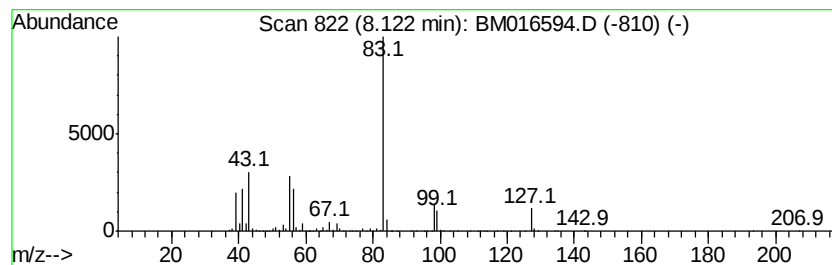
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 1-Butene, 2,3,3-trimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.12	163.80 ng	6603340	1,4-Dichlorobenzene-d4	8.01

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Butene, 2,3,3-trimethyl-	98	C7H14	000594-56-9	72
2		1H-Imidazol-2-amine	83	C3H5N3	007720-39-0	53
3		1,5-Heptadien-4-one, 3,3,6-trime...	152	C10H16O	000546-49-6	47
4		1H-Pyrazole, 4,5-dihydro-5,5-dim...	98	C5H10N2	004320-85-8	40
5		2-Pentene, 4,4-dimethyl-, (E)-	98	C7H14	000690-08-4	40



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM082418\
Data File : BM016594.D
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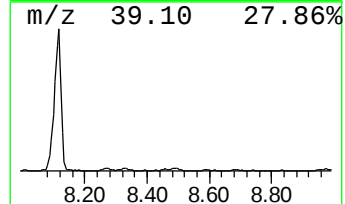
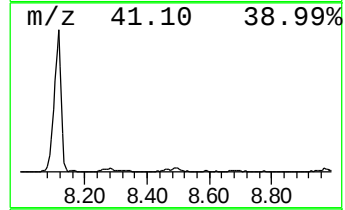
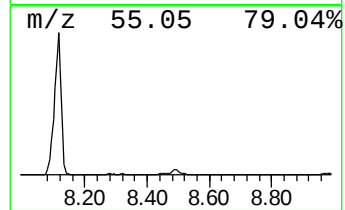
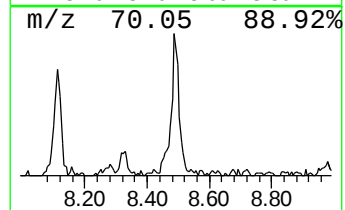
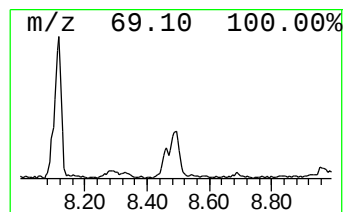
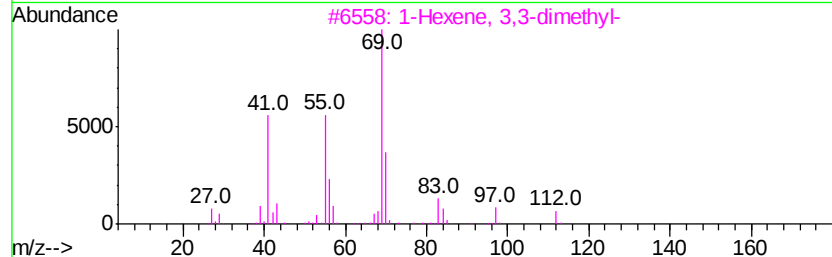
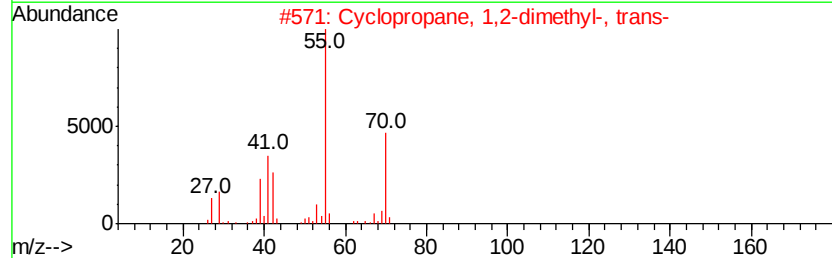
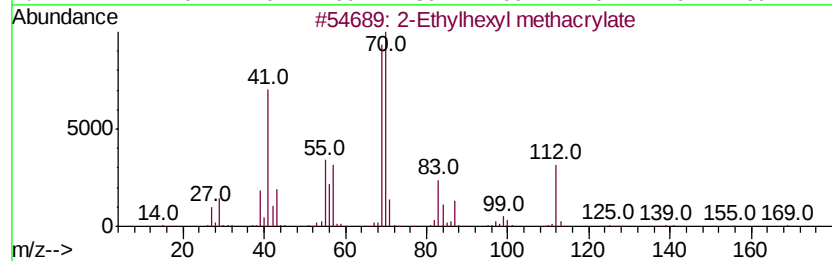
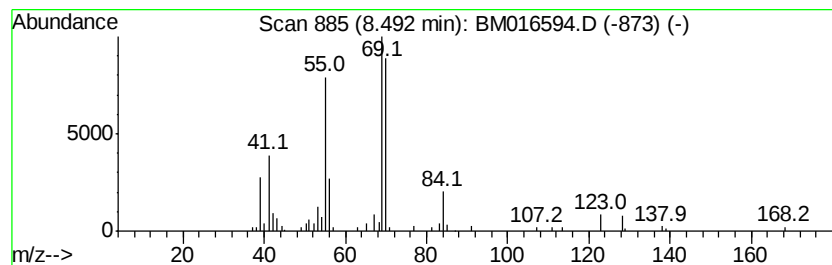
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 2-Ethylhexyl methacrylate Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.	
8.49	6.23 ng	251340	1,4-Dichlorobenzene-d4	8.01	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Ethylhexyl methacrylate	198	C12H22O2	000688-84-6	59
2	Cyclopropane, 1,2-dimethyl-, trans-	70	C5H10	002402-06-4	38
3	1-Hexene, 3,3-dimethyl-	112	C8H16	003404-77-1	38
4	Cyclohexane, 1,2,3-trimethyl-	126	C9H18	001678-97-3	36
5	Cyclopropane, 1,2-dimethyl-, cis-	70	C5H10	000930-18-7	32



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM082418\
Data File : BM016594.D
Acq On : 24 Aug 2018 18:39
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Sample : J4616-01
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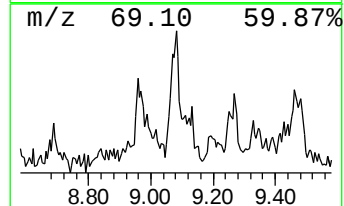
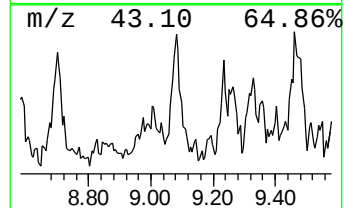
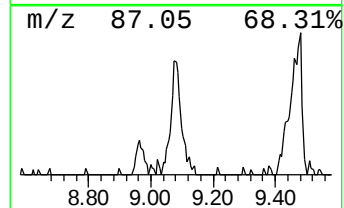
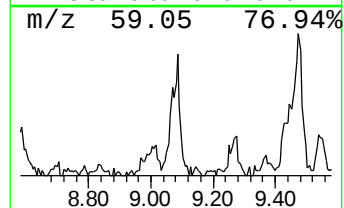
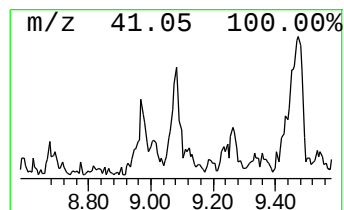
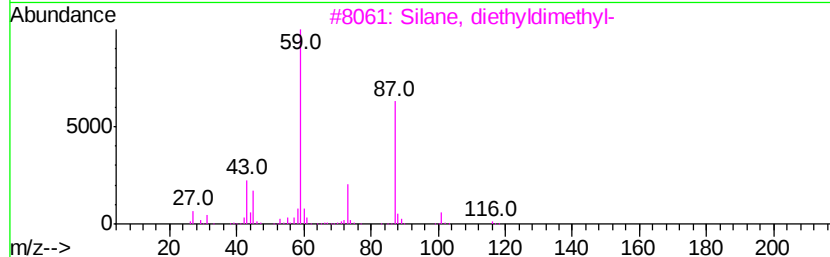
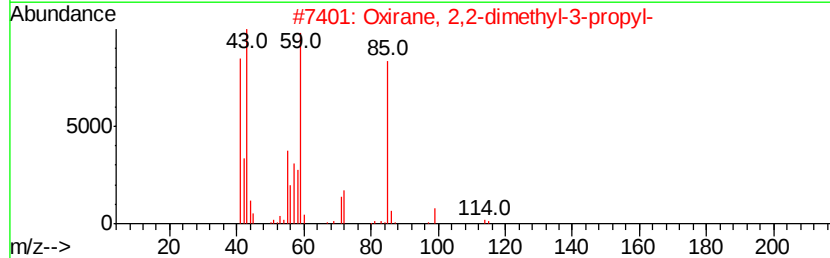
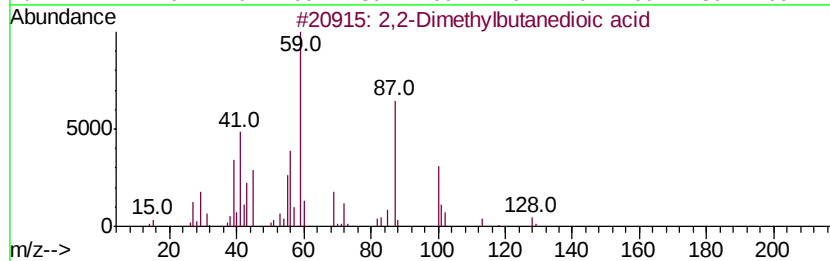
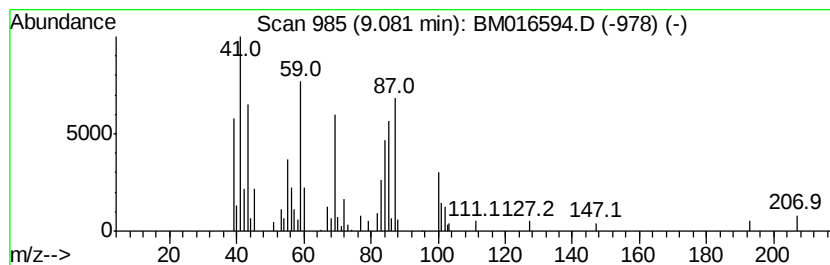
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 unknown9.08 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.08	4.34 ng	174926	1,4-Dichlorobenzene-d4	8.01

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2,2-Dimethylbutanedioic acid	146	C6H10O4	000597-43-3	47
2			Oxirane, 2,2-dimethyl-3-propyl-	114	C7H14O	017612-35-0	35
3			Silane, diethyldimethyl-	116	C6H16Si	000756-81-0	35
4			3-Octanol	130	C8H18O	000589-98-0	27
5			2-Propanol, 1-(1,3-dimethylbutoxy)-	160	C9H20O2	054340-89-5	27



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM082418\
Data File : BM016594.D
Acq On : 24 Aug 2018 18:39
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Misc :
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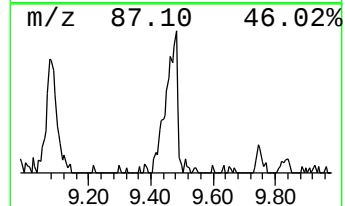
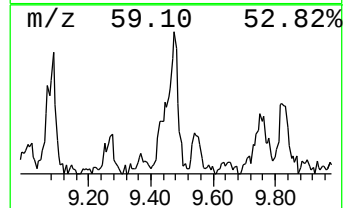
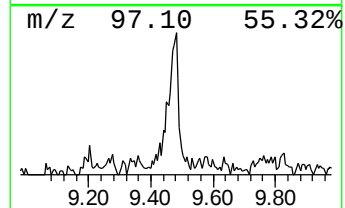
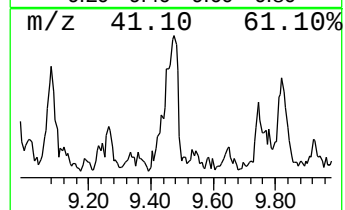
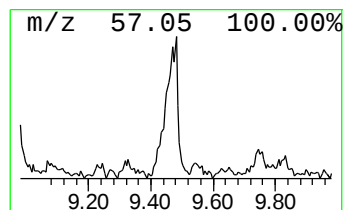
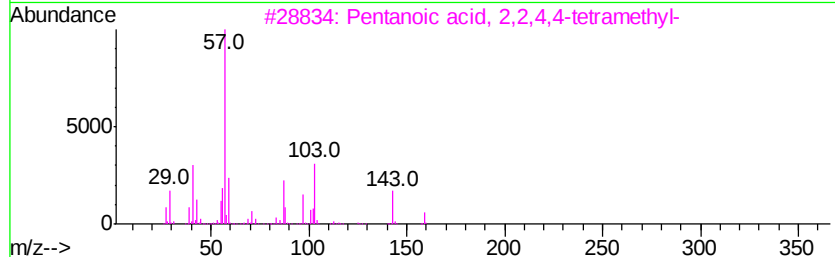
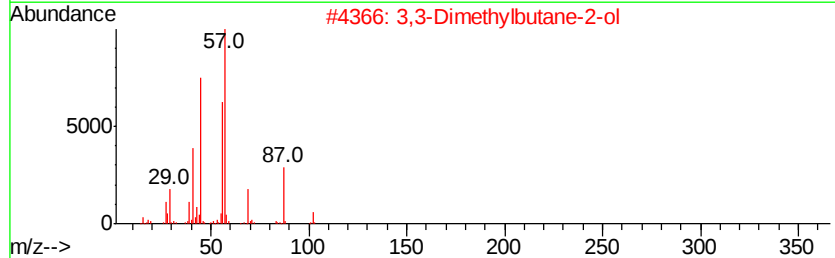
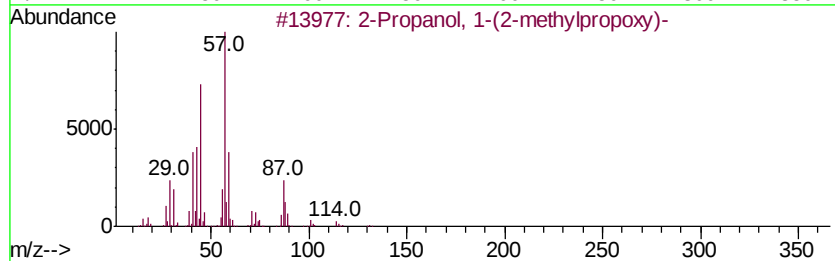
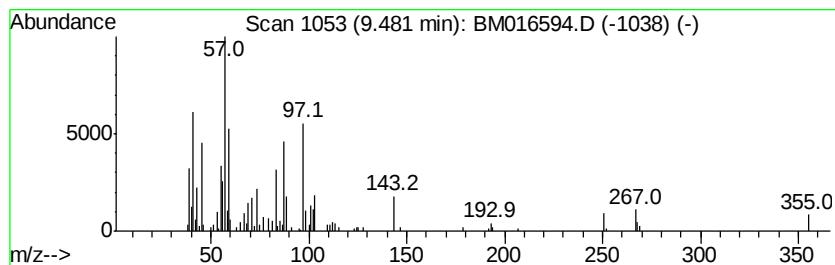
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 unknown9.48 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.48	8.76 ng	457121	Naphthalene-d8	10.81

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanol, 1-(2-methylpropoxy)-	132	C7H16O2	023436-19-3	38		
2	3,3-Dimethylbutane-2-ol	102	C6H14O	000464-07-3	22		
3	Pentanoic acid, 2,2,4,4-tetramet...	158	C9H18O2	003302-12-3	16		
4	Propane, 1-methoxy-2,2-dimethyl-	102	C6H14O	001118-00-9	14		
5	3-Pentanol, 2,2-dimethyl-	116	C7H16O	003970-62-5	12		



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM082418\
Data File : BM016594.D
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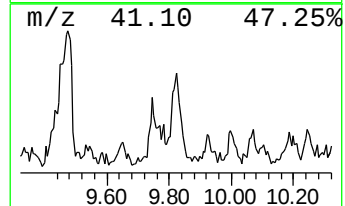
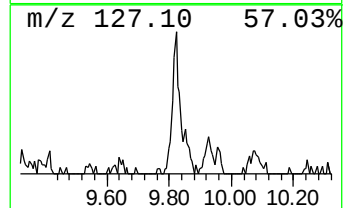
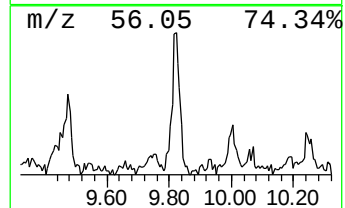
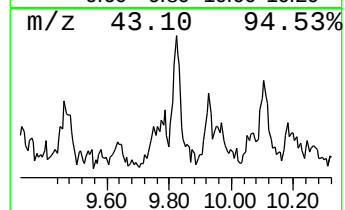
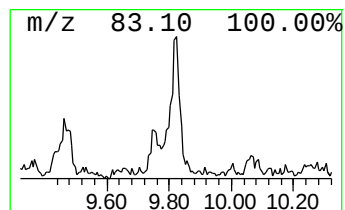
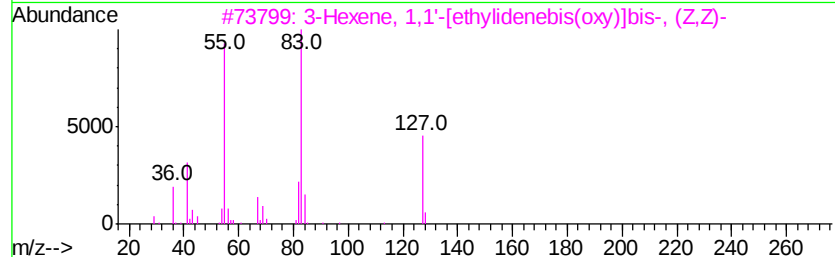
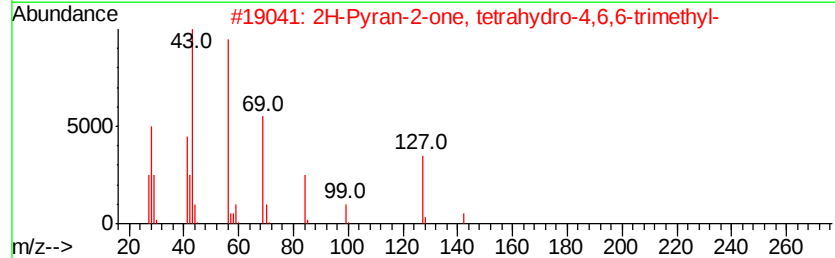
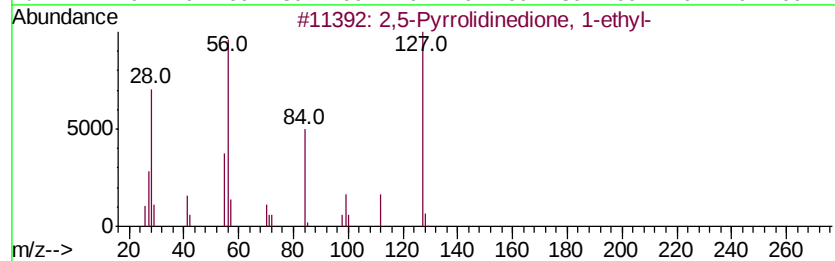
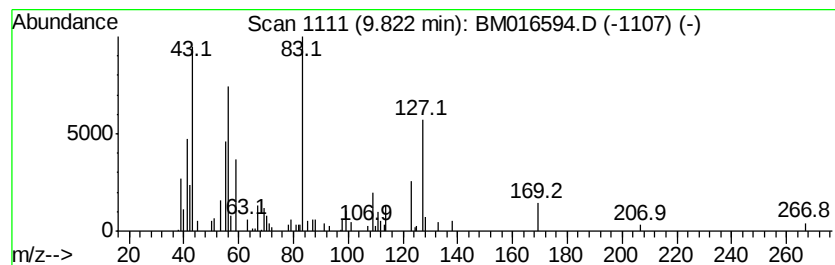
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 unknown9.82 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.82	4.82 ng	251791	Naphthalene-d8	10.81

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,5-Pyrrolidinedione, 1-ethyl-	127	C6H9NO2	002314-78-5	30		
2	2H-Pyran-2-one, tetrahydro-4,6,6...	142	C8H14O2	020628-36-8	25		
3	3-Hexene, 1,1'-[ethylidenebis(ox...	226	C14H26O2	063449-64-9	16		
4	Cyclobutane, 1,2-diethyl-	112	C8H16	061141-83-1	14		
5	4-Piperidinemethanamine	114	C6H14N2	007144-05-0	14		



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM082418\
Data File : BM016594.D
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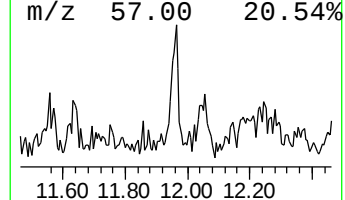
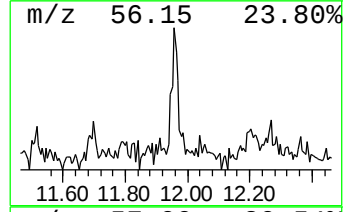
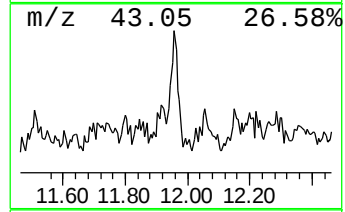
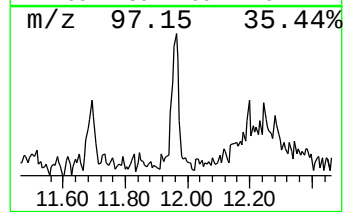
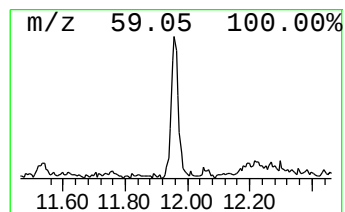
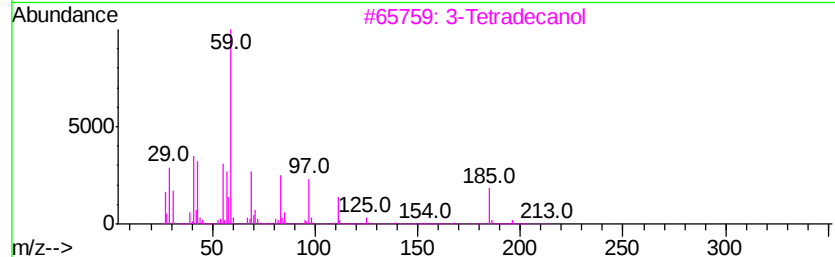
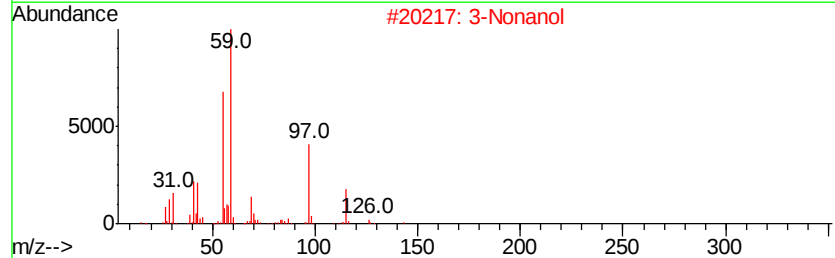
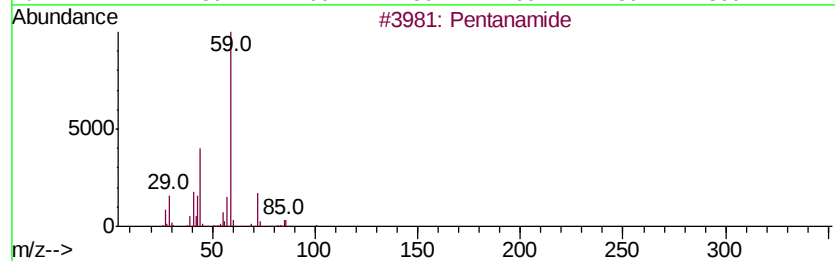
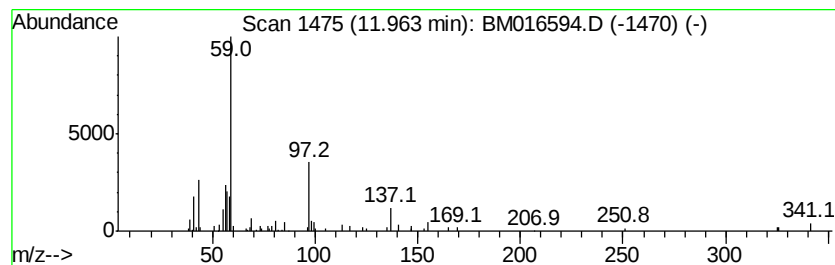
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 unknown11.96 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.96	4.09 ng	213744	Naphthalene-d8	10.81

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentanamide	101	C5H11NO	000626-97-1	43
2			3-Nonanol	144	C9H20O	000624-51-1	40
3			3-Tetradecanol	214	C14H30O	001653-32-3	40
4			4-Pentene-2-ol, 2-methyl	100	C6H12O	000624-97-5	38
5			Hexylene Glycol	118	C6H14O2	000107-41-5	38



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM082418\
Data File : BM016594.D
Acq On : 24 Aug 2018 18:39
Operator : SJ/JU
Sample : J4616-01
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
VB27184

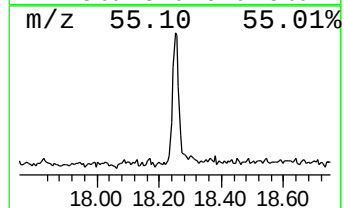
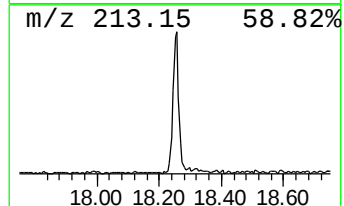
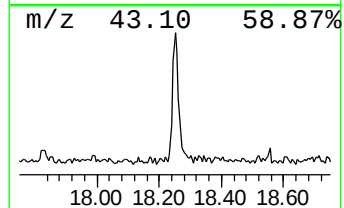
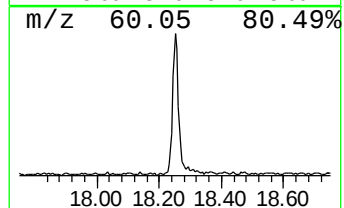
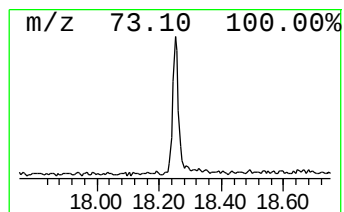
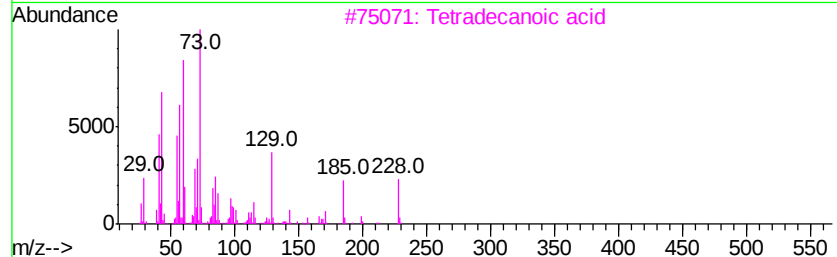
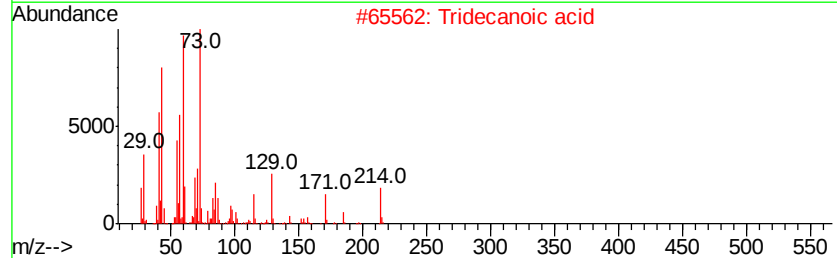
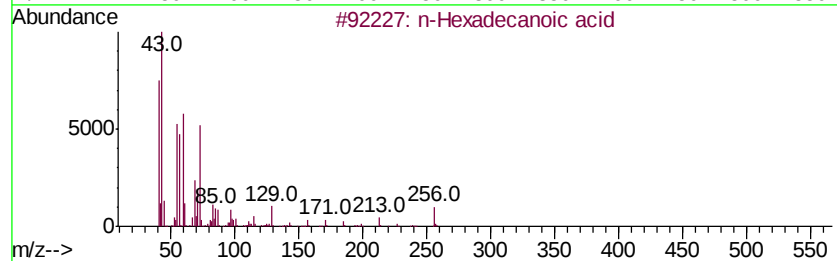
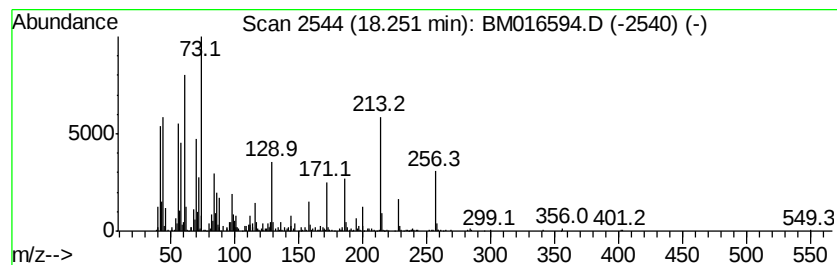
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.25	11.34 ng	1410330	Phenanthrene-d10	17.39

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
2		Tridecanoic acid	214	C13H26O2	000638-53-9	91
3		Tetradecanoic acid	228	C14H28O2	000544-63-8	52
4		Dodecanoic acid	200	C12H24O2	000143-07-7	38
5		Nonanoic acid	158	C9H18O2	000112-05-0	38



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM082418\
Data File : BM016594.D
Acq On : 24 Aug 2018 18:39
Operator : SJ/JU
Sample : J4616-01
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
VB27184

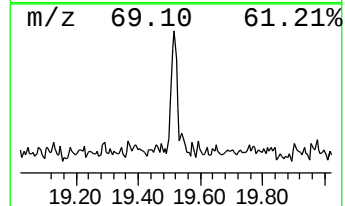
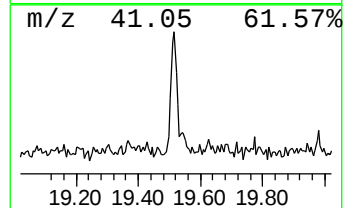
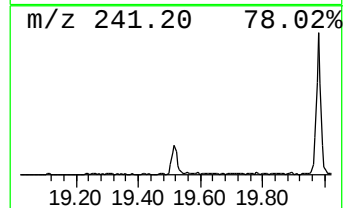
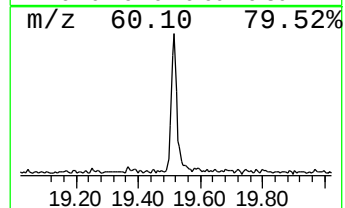
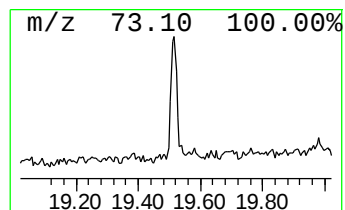
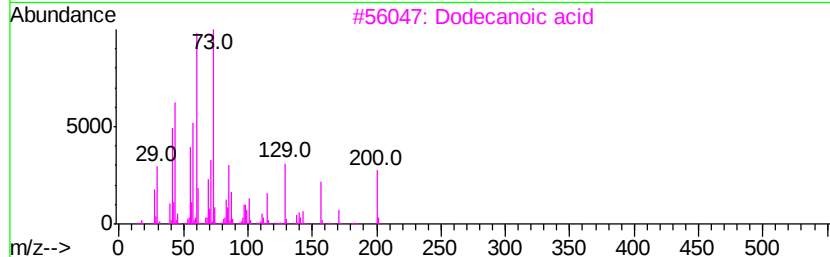
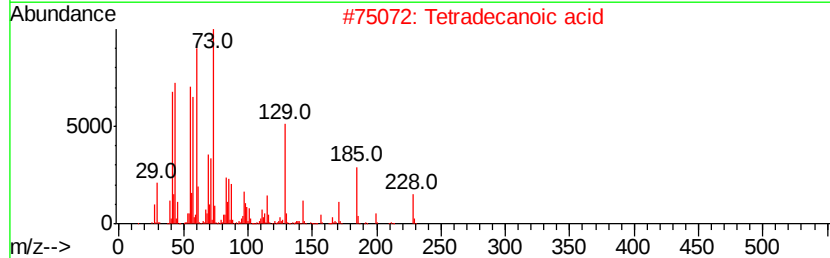
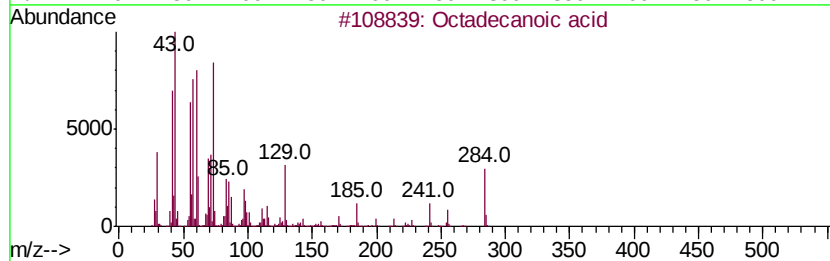
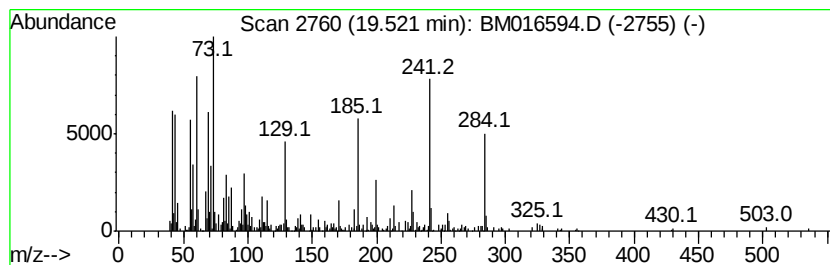
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Octadecanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.52	4.62 ng	792196	Chrysene-d12	21.53

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	90
2			Tetradecanoic acid	228	C14H28O2	000544-63-8	43
3			Dodecanoic acid	200	C12H24O2	000143-07-7	38
4			Undecanoic acid	186	C11H22O2	000112-37-8	38
5			Tridecanoic acid	214	C13H26O2	000638-53-9	30



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM082418\
Data File : BM016594.D
Acq On : 24 Aug 2018 18:39
Operator : SJ/JU
Sample : J4616-01
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
VB27184

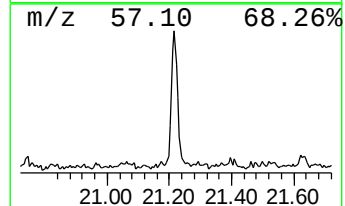
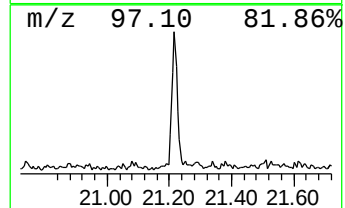
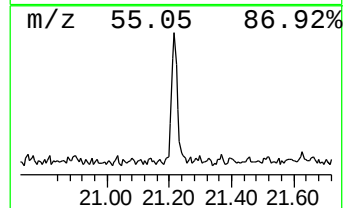
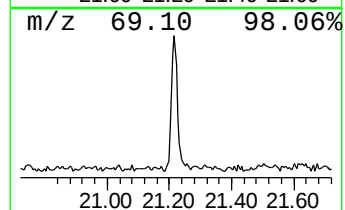
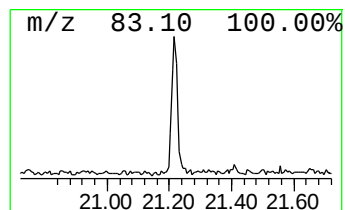
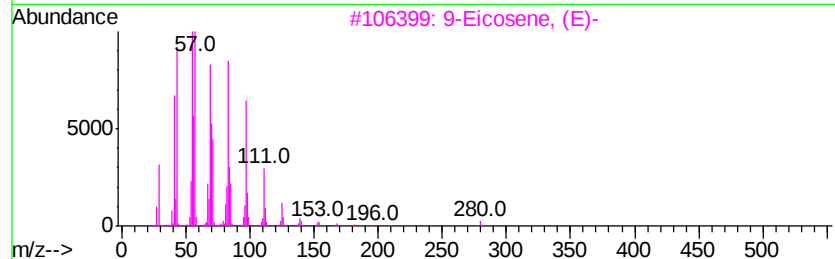
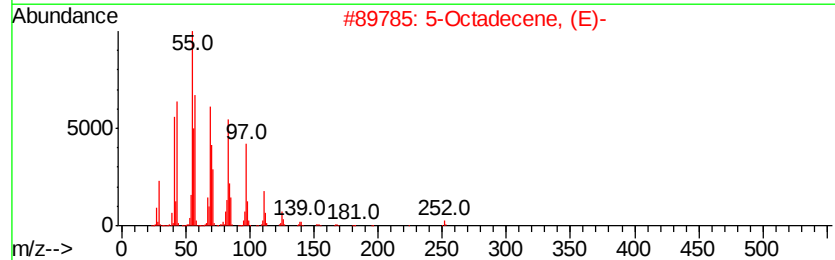
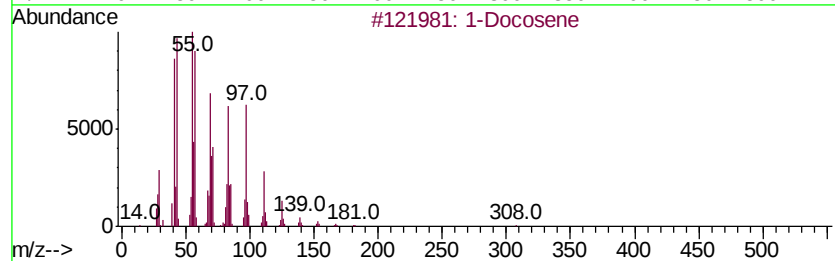
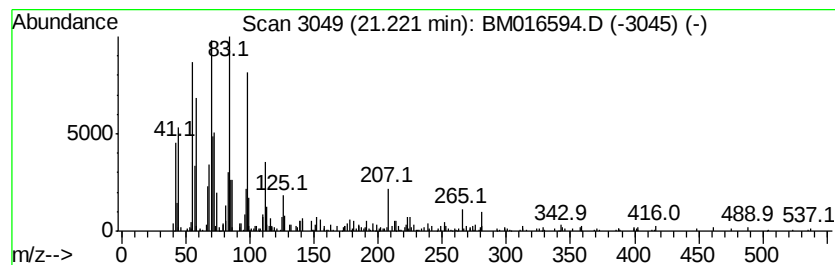
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM082118.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 1-Docosene Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.22	3.86 ng	663256	Chrysene-d12	21.53

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Docosene	308	C22H44	001599-67-3	93
2			5-Octadecene, (E)-	252	C18H36	007206-21-5	91
3			9-Eicosene, (E)-	280	C20H40	074685-29-3	91
4			3-Eicosene, (E)-	280	C20H40	074685-33-9	91
5			2-Tetradecene, (E)-	196	C14H28	035953-53-8	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM082418\
 Data File : BM016594.D
 Acq On : 24 Aug 2018 18:39
 Operator : SJ/JU
 Sample : J4616-01
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 VB27184

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\8270-BM082118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanol, 2-met...	3.62	5.2	ng	210297	1	8.01	806271	20.0
3-Octanol	4.35	4.1	ng	164069	1	8.01	806271	20.0
Butanamide, 3,3-d...	5.39	4.0	ng	160420	1	8.01	806271	20.0
unknown7.12	7.12	8.9	ng	357691	1	8.01	806271	20.0
unknown7.53	7.53	70.3	ng	2831900	1	8.01	806271	20.0
2-Pentene, 4-meth...	7.88	7.5	ng	300557	1	8.01	806271	20.0
1-Butene, 2,3,3-t...	8.12	163.8	ng	6603340	1	8.01	806271	20.0
2-Ethylhexyl meth...	8.49	6.2	ng	251340	1	8.01	806271	20.0
unknown9.08	9.08	4.3	ng	174926	1	8.01	806271	20.0
unknown9.48	9.48	8.8	ng	457121	2	10.81	1044020	20.0
unknown9.82	9.82	4.8	ng	251791	2	10.81	1044020	20.0
unknown11.96	11.96	4.1	ng	213744	2	10.81	1044020	20.0
n-Hexadecanoic acid	18.25	11.3	ng	1410330	4	17.39	2487380	20.0
Octadecanoic acid	19.52	4.6	ng	792196	5	21.53	3432290	20.0
1-Docosene	21.22	3.9	ng	663256	5	21.53	3432290	20.0