

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM071118\
 Data File : BM015887.D
 Acq On : 11 Jul 2018 14:01
 Operator : SJ/JU
 Sample : J3925-02
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 MW-2

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-BM070518.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	3.816	84	90	99	rBV	208535	279770	5.03%	0.846%
2	5.316	340	345	357	rVB	75621	129824	2.34%	0.393%
3	5.798	421	427	442	rBV	695019	1124677	20.23%	3.402%
4	7.445	701	707	719	rBV	414268	725022	13.04%	2.193%
5	7.816	763	770	787	rBV	1608196	2621006	47.14%	7.927%
6	8.292	845	851	858	rBV	376544	615061	11.06%	1.860%
7	8.522	881	890	899	rBV3	34939	72551	1.30%	0.219%
8	8.616	899	906	922	rVB	1595367	2623431	47.19%	7.934%
9	9.328	1020	1027	1030	rBV2	43297	74594	1.34%	0.226%
10	9.480	1046	1053	1060	rBV	1251553	2074008	37.30%	6.273%
11	10.116	1156	1161	1175	rVB5	51065	128907	2.32%	0.390%
12	11.122	1324	1332	1344	rBV	452607	830861	14.94%	2.513%
13	11.398	1368	1379	1386	rBV2	23507	62716	1.13%	0.190%
14	12.080	1491	1495	1499	rBV2	33763	60104	1.08%	0.182%
15	12.363	1534	1543	1550	rBV2	50998	110335	1.98%	0.334%
16	13.451	1724	1728	1733	rBV2	40952	63191	1.14%	0.191%
17	13.539	1736	1743	1749	rBV	3130604	4372660	78.65%	13.225%
18	13.751	1773	1779	1789	rBV	1124837	1635145	29.41%	4.945%
19	14.363	1878	1883	1888	rVB	141560	204297	3.67%	0.618%
20	14.539	1907	1913	1919	rBV6	47726	85433	1.54%	0.258%
21	14.915	1971	1977	1985	rVB2	629801	978796	17.61%	2.960%
22	15.539	2075	2083	2090	rVB4	38565	113099	2.03%	0.342%
23	15.633	2095	2099	2108	rVB	98116	158842	2.86%	0.480%
24	16.392	2222	2228	2240	rBV	2496968	3561885	64.07%	10.773%
25	17.639	2434	2440	2448	rBV	752039	1122666	20.19%	3.395%
26	18.268	2544	2547	2552	rVB3	70623	80344	1.45%	0.243%
27	18.480	2578	2583	2593	rBV	276775	400706	7.21%	1.212%
28	18.562	2593	2597	2604	rVB	59918	89784	1.61%	0.272%
29	18.692	2616	2619	2628	rVB	45709	74182	1.33%	0.224%
30	19.309	2720	2724	2732	rVB2	45215	77793	1.40%	0.235%
31	19.615	2773	2776	2782	rVB3	45975	74462	1.34%	0.225%
32	19.727	2791	2795	2807	rBV	154108	224157	4.03%	0.678%
33	20.021	2842	2845	2851	rVB3	37793	55755	1.00%	0.169%
34	20.203	2870	2876	2887	rVB	4731579	5559613	100.00%	16.815%

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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	21.503	3093	3097	3104	rBV	312947	381143	6.86%	1.153%
36	21.750	3134	3139	3144	rBV	923856	1177159	21.17%	3.560%
37	24.144	3540	3546	3555	rVB	532572	1039681	18.70%	3.144%

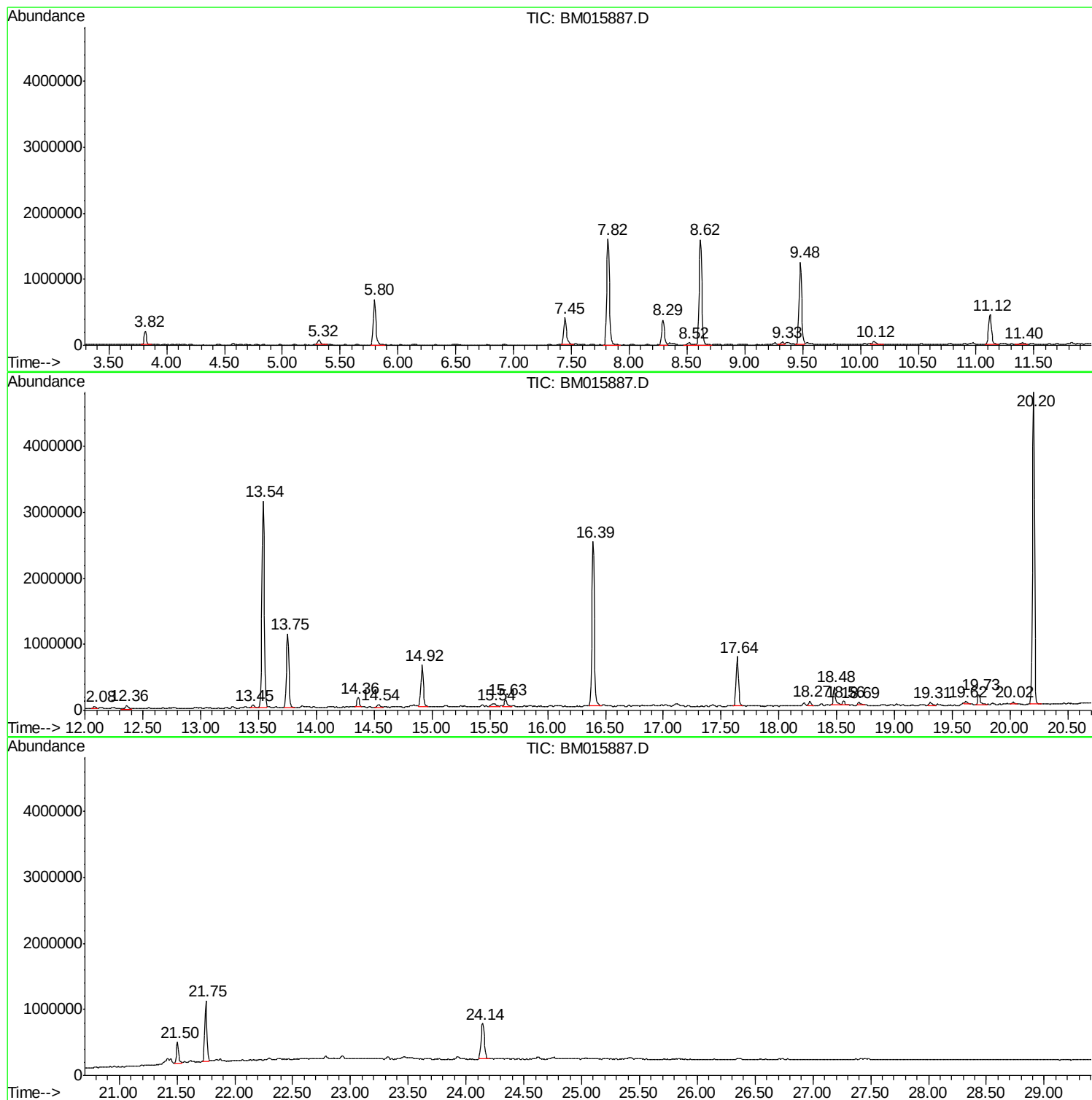
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM070518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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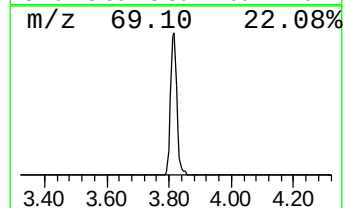
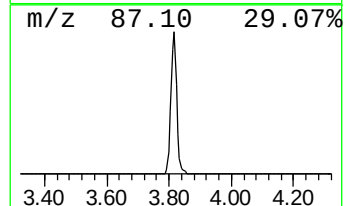
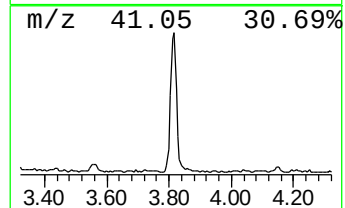
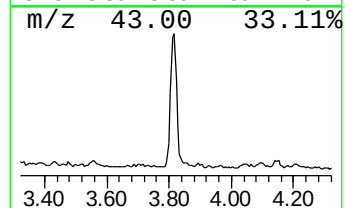
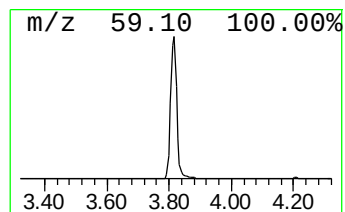
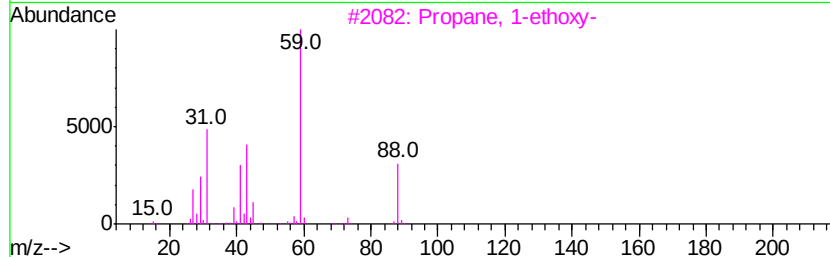
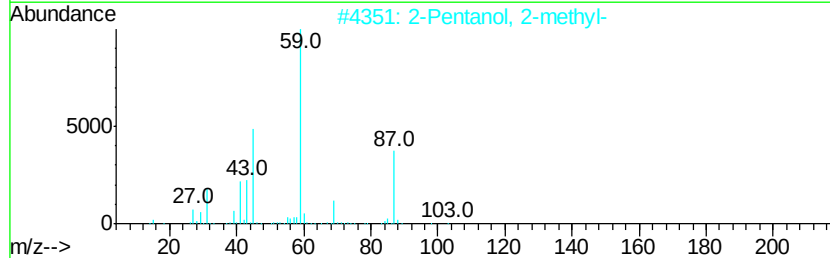
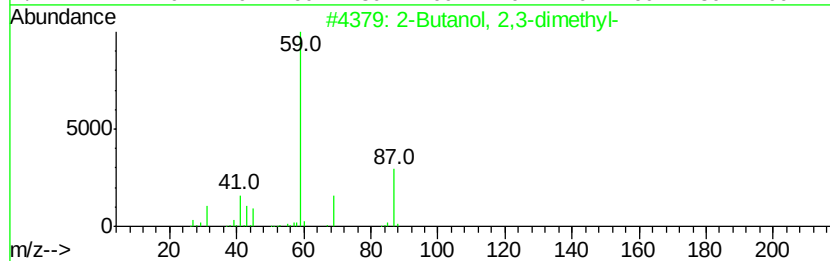
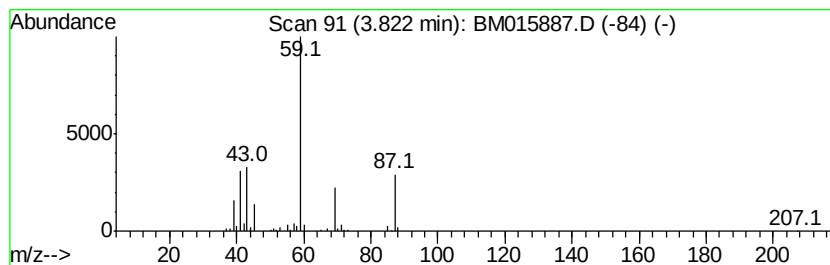
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Butanol, 2,3-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.82	9.10 ng	279770	1,4-Dichlorobenzene-d4	8.29

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Butanol, 2,3-dimethyl-	102	C6H14O	000594-60-5	78
2			2-Pentanol, 2-methyl-	102	C6H14O	000590-36-3	72
3			Propane, 1-ethoxy-	88	C5H12O	000628-32-0	59
4			Propane, 2-ethoxy-2-methyl-	102	C6H14O	000637-92-3	50
5			Amylene Hydrate	88	C5H12O	000075-85-4	45



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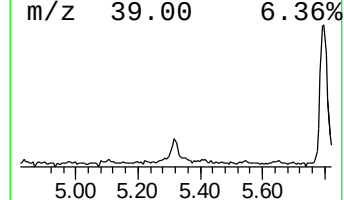
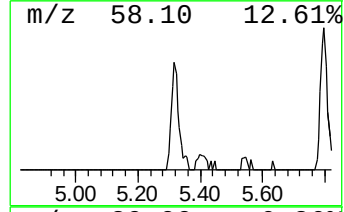
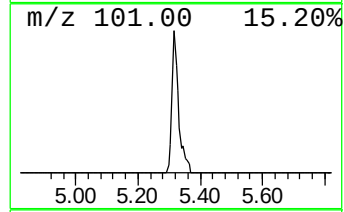
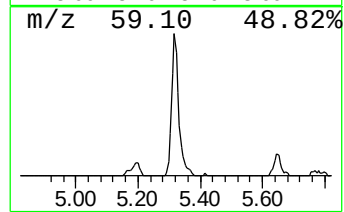
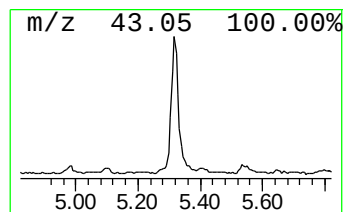
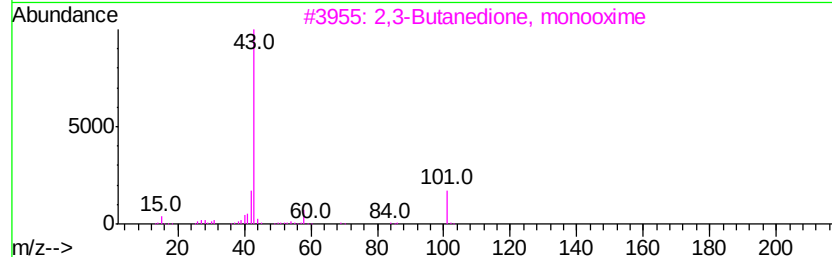
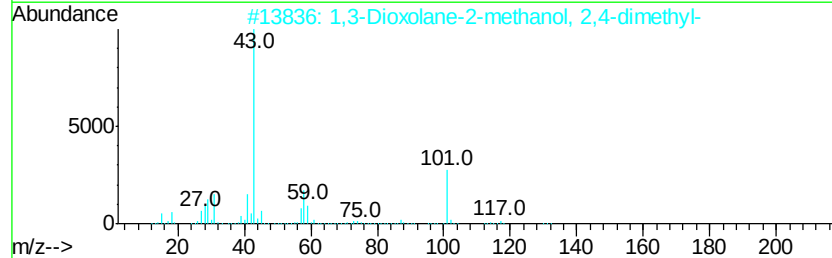
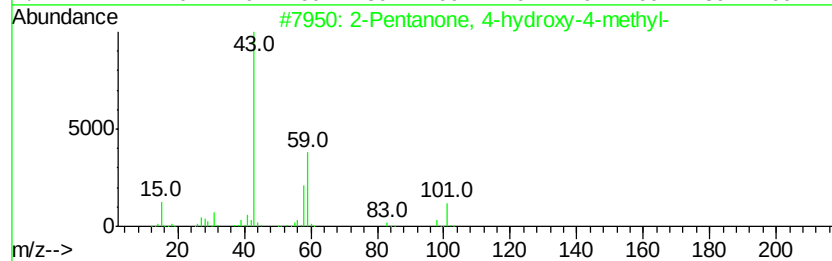
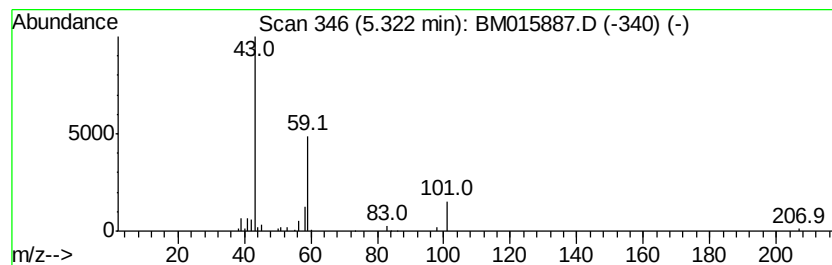
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.32	4.22 ng	129824	1,4-Dichlorobenzene-d4	8.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2		1,3-Dioxolane-2-methanol, 2,4-di...	132	C6H12O3	053951-43-2	17
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16
4		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5		4-Penten-2-one, 4-methyl-	98	C6H10O	003744-02-3	9



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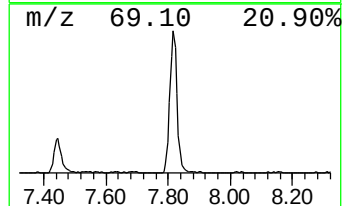
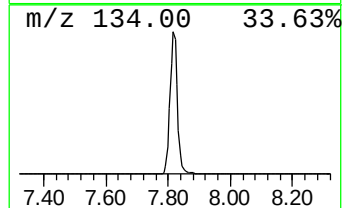
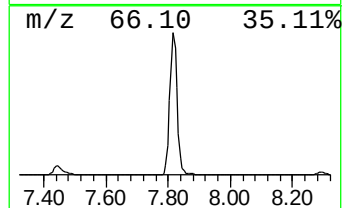
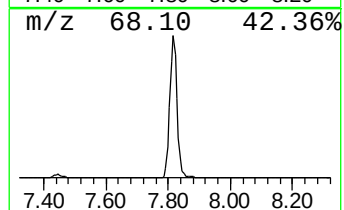
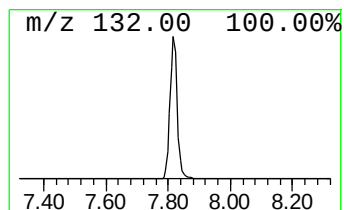
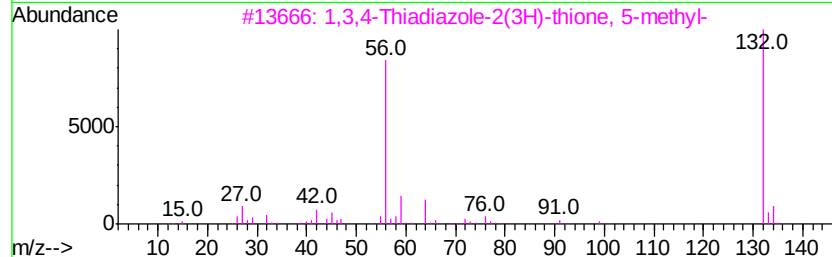
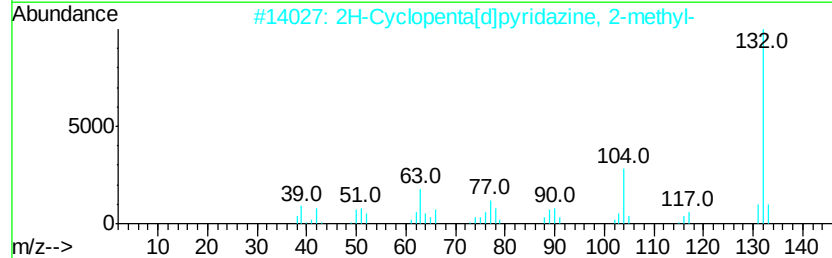
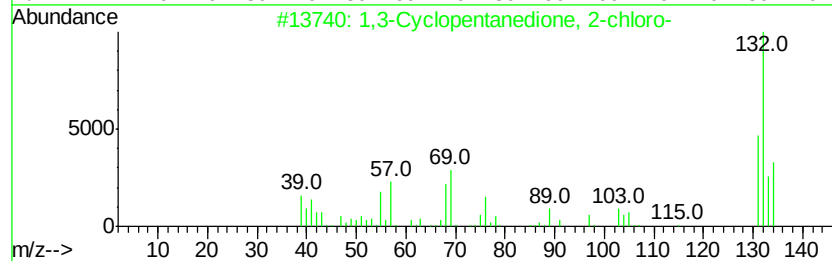
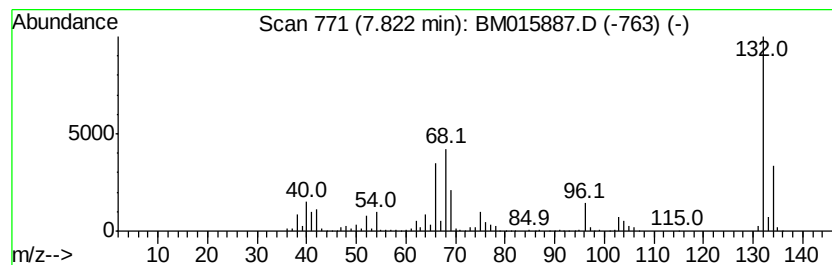
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown7.82 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.82	85.23 ng	2621010	1,4-Dichlorobenzene-d4	8.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	35
2		2H-Cyclopenta[d]pyridazine, 2-methyl-	132	C8H8N2	022291-85-6	27
3		1,3,4-Thiadiazole-2(3H)-thione, 5-methyl-	132	C3H4N2S2	029490-19-5	22
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	22
5		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22



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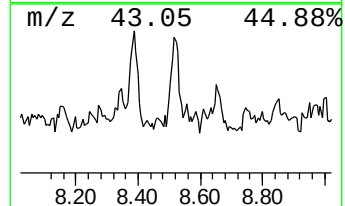
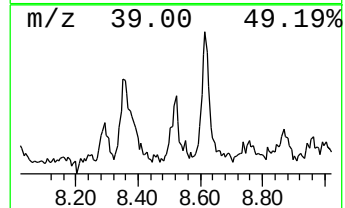
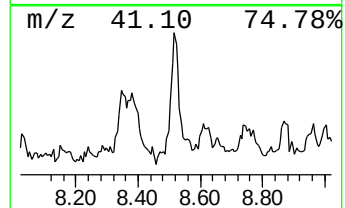
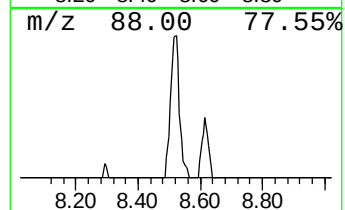
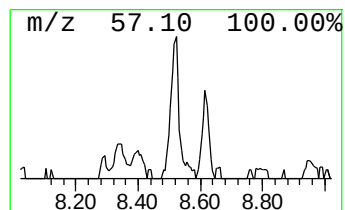
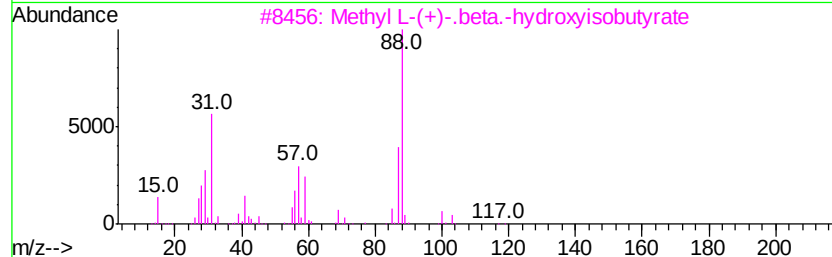
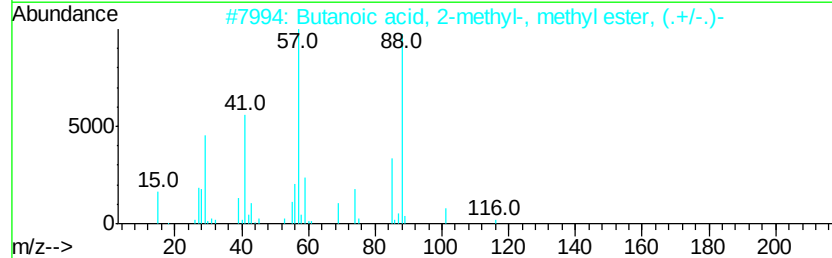
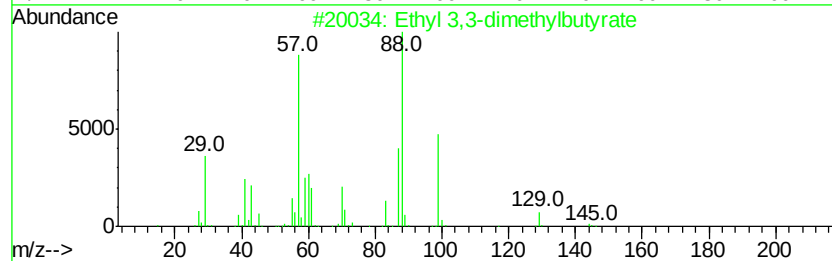
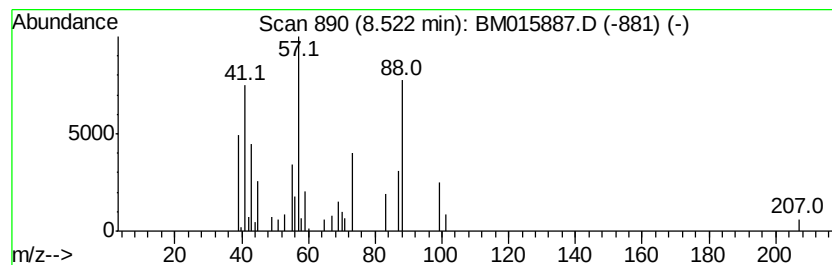
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Ethyl 3,3-dimethylbutyrate Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.52	2.36 ng	72551	1,4-Dichlorobenzene-d4	8.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Ethyl 3,3-dimethylbutyrate	144	C8H16O2	005340-78-3	50
2		Butanoic acid, 2-methyl-, methyl...	116	C6H12O2	053955-81-0	38
3		Methyl L-(+)-.beta.-hydroxyisobut...	118	C5H10O3	080657-57-4	32
4		Methyl (R)-(-)-3-hydroxy-2-methyl...	118	C5H10O3	072657-23-9	32
5		Butanoic acid, 2-methyl-, methyl...	116	C6H12O2	000868-57-5	32



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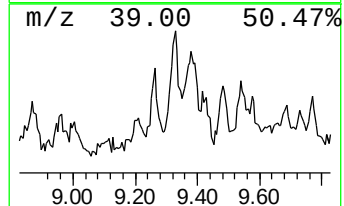
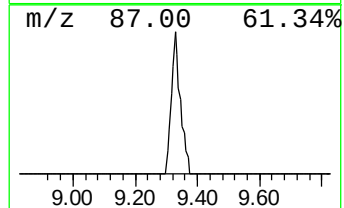
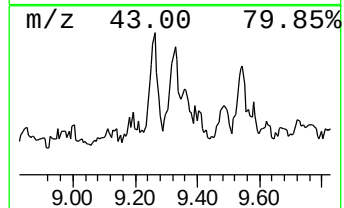
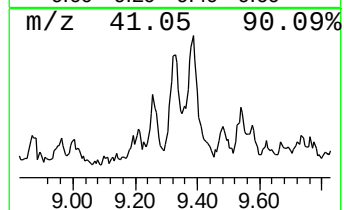
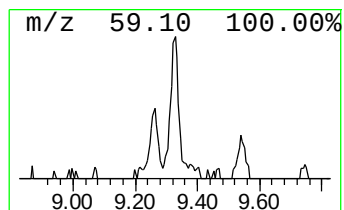
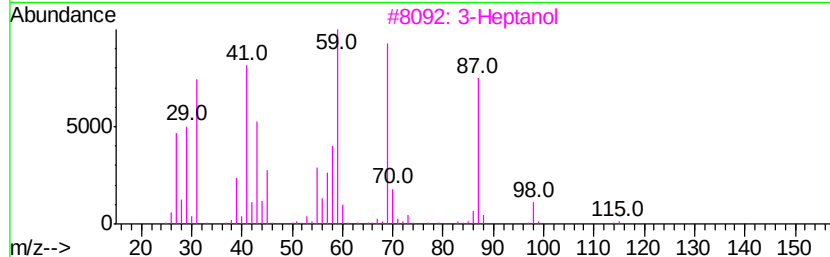
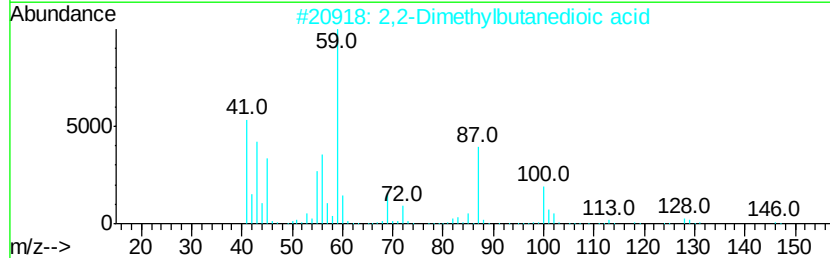
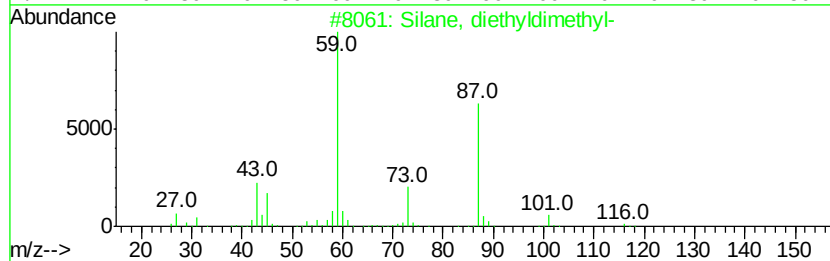
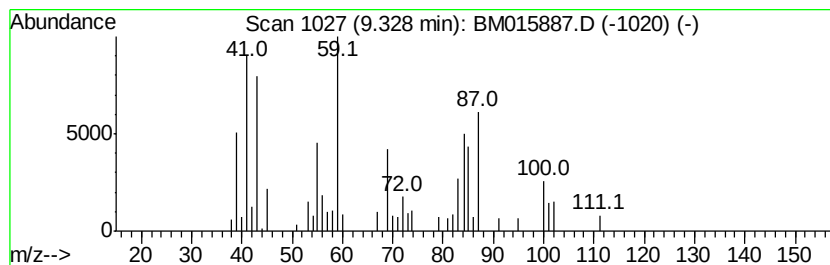
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Peak Number 6 unknown9.33 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.33	2.43 ng	74594	1,4-Dichlorobenzene-d4	8.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Silane, diethyldimethyl-	116	C6H16Si	000756-81-0	38
2		2,2-Dimethylbutanedioic acid	146	C6H10O4	000597-43-3	38
3		3-Heptanol	116	C7H16O	000589-82-2	25
4		1,3-Propanediol, 2,2-diethyl-	132	C7H16O2	000115-76-4	25
5		2-Pentene, 2-methoxy-	100	C6H12O	061142-47-0	18



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM071118\
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Acq On : 11 Jul 2018 14:01
Operator : SJ/JU
Sample : J3925-02
Misc :
ALS Vial : 4 Sample Multiplier: 1

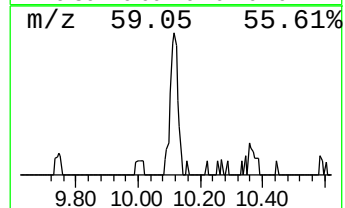
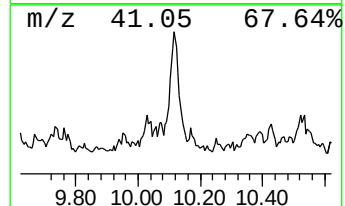
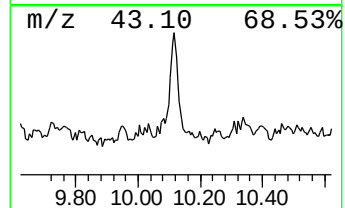
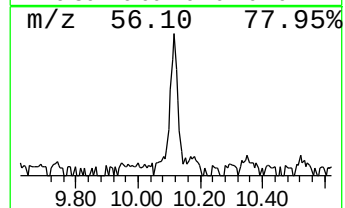
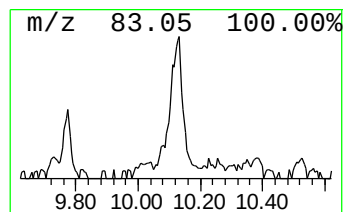
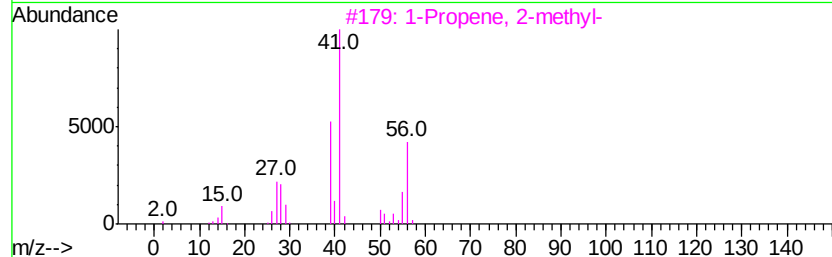
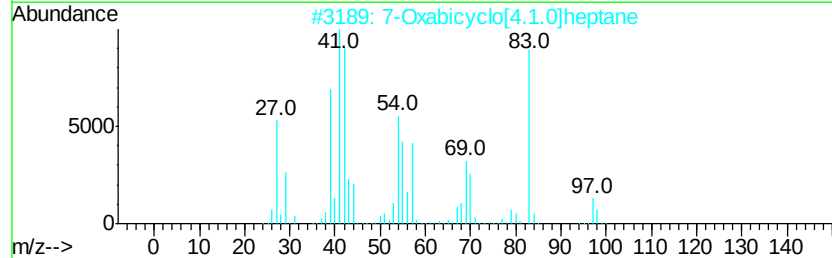
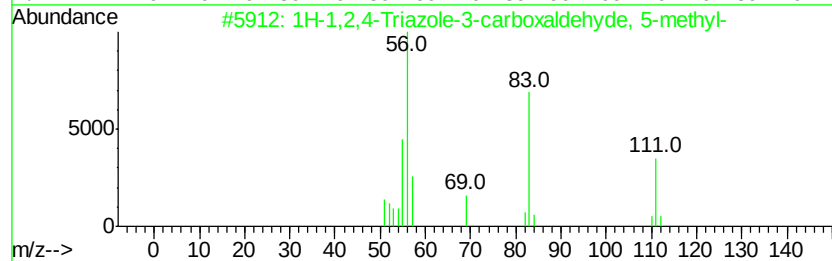
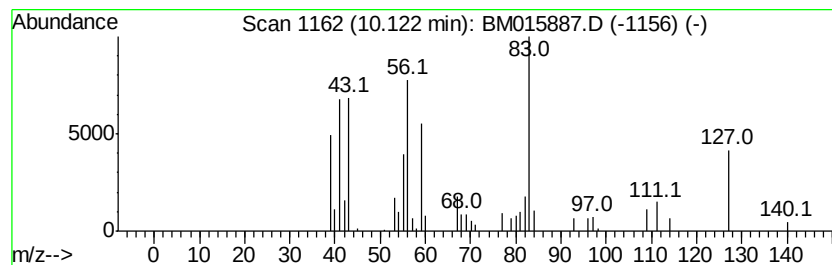
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ClientSampled :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM070518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 1H-1,2,4-Triazole-3-carboxal... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.12	3.10 ng	128907	Naphthalene-d8	11.12	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-1,2,4-Triazole-3-carboxaldehv...	111	C4H5N3O	056804-98-9	50
2	7-Oxabicyclo[4.1.0]heptane	98	C6H10O	000286-20-4	35
3	1-Propene, 2-methyl-	56	C4H8	000115-11-7	30
4	6-Methyl-2-heptyl methylphosphon...	210	C9H20FO2P	159395-76-3	27
5	Urea, N,N'-di-2-propenyl-	140	C7H12N2O	001801-72-5	18



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM071118\
Data File : BM015887.D
Acq On : 11 Jul 2018 14:01
Operator : SJ/JU
Sample : J3925-02
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ALS Vial : 4 Sample Multiplier: 1

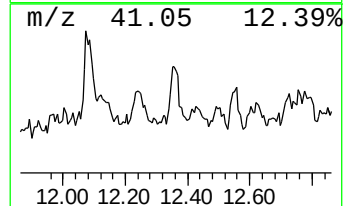
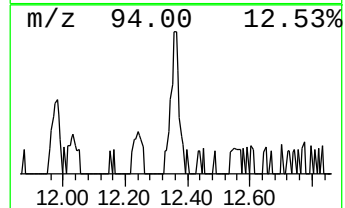
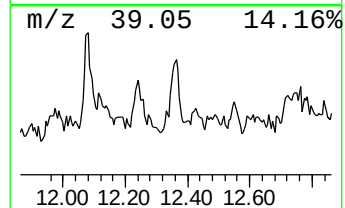
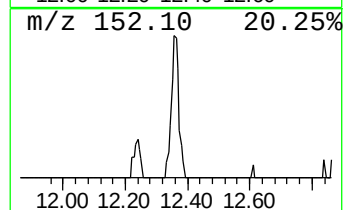
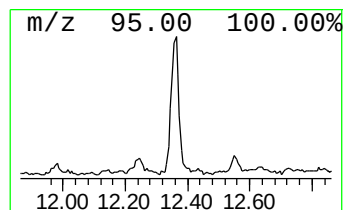
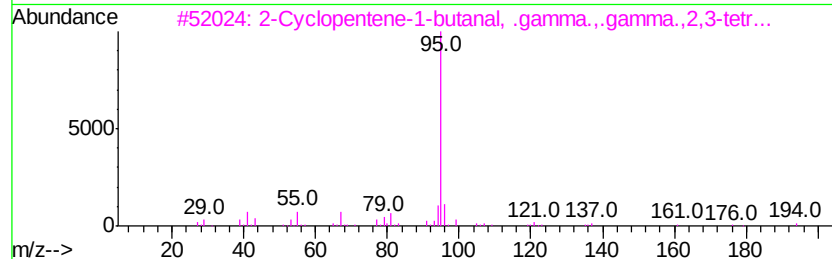
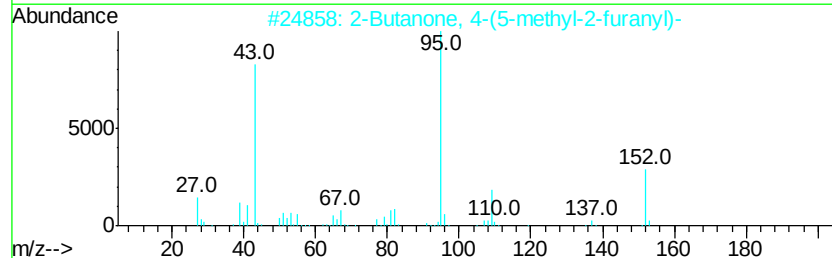
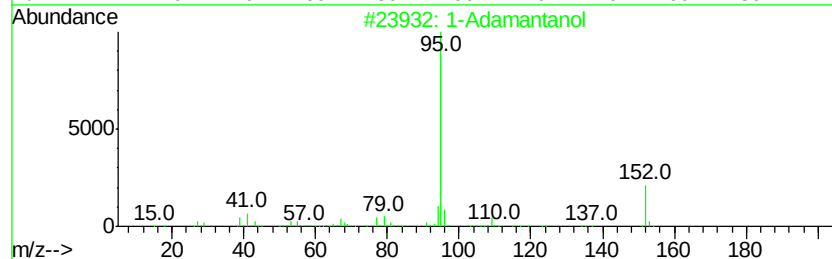
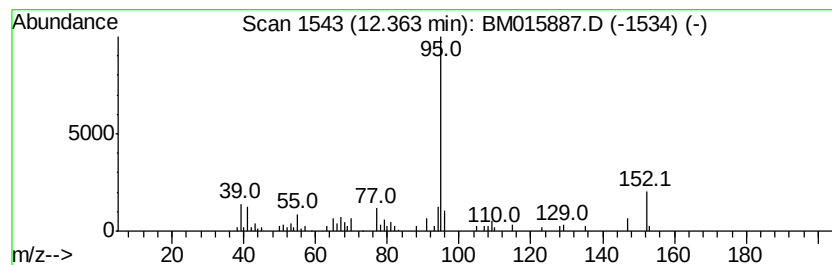
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 1-Adamantanol Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.36	2.66 ng	110335	Naphthalene-d8	11.12
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1-Adamantanol	152 C10H16O	000768-95-6	87
2	2-Butanone, 4-(5-methyl-2-furanyl)-	152 C9H12O2	013679-56-6	59
3	2-Cyclopentene-1-butanal, .gamma...	194 C13H22O	060714-25-2	50
4	4-Pyridinol	95 C5H5NO	000626-64-2	49
5	2-Norbornyl bromide	174 C7H11Br	029342-65-2	47



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM071118\
Data File : BM015887.D
Acq On : 11 Jul 2018 14:01
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Sample : J3925-02
Misc :
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Instrument :
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ClientSampleId :
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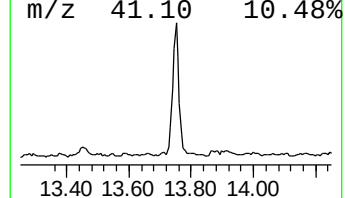
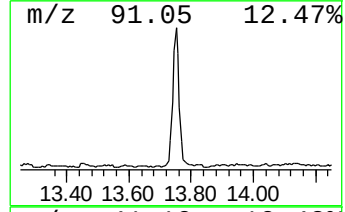
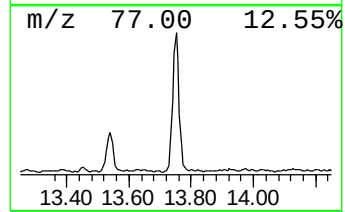
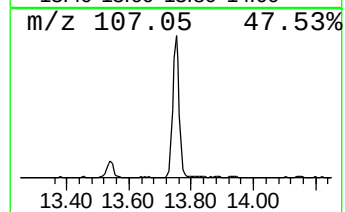
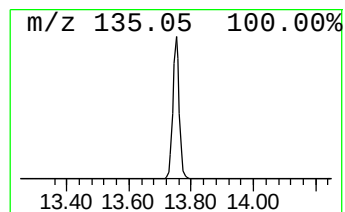
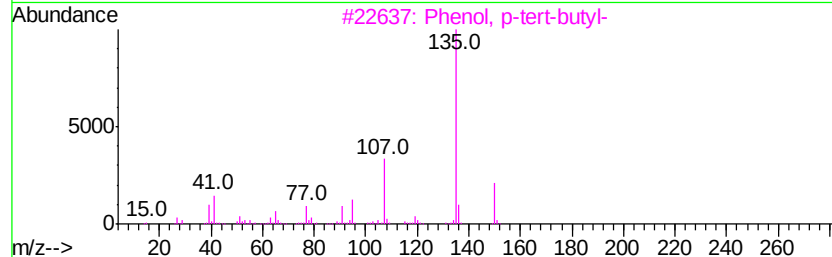
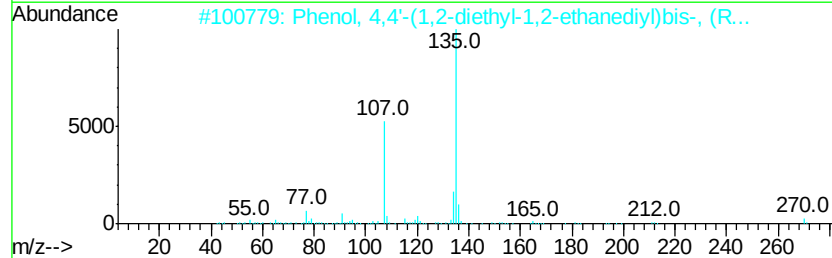
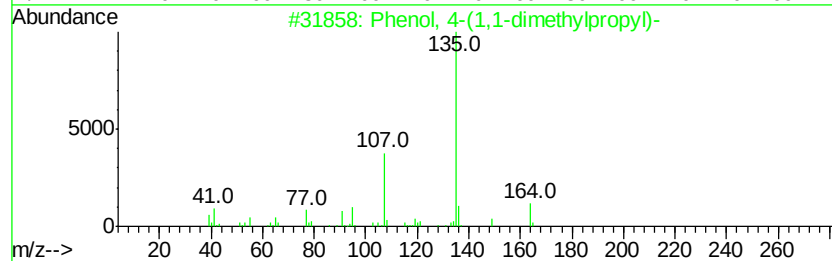
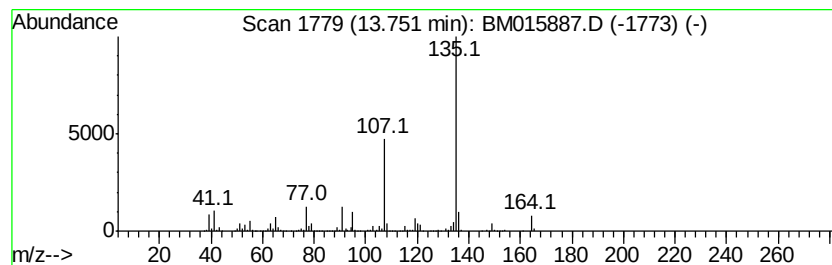
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Phenol, 4-(1,1-dimethylprop... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.75	33.41 ng	1635150	Acenaphthene-d10	14.92

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	95
2		Phenol, 4,4'-(1,2-diethyl-1,2-eth...	270	C18H22O2	000084-16-2	72
3		Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	72
4		4,6-Bis(4-ethoxybenzylthio)-5-ni...	457	C22H23N3O4S2	325957-76-4	64
5		2-Aziridinone, 1-(1-adamantyl)-3...	273	C18H27NO	026905-18-0	53



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM071118\
Data File : BM015887.D
Acq On : 11 Jul 2018 14:01
Operator : SJ/JU
Sample : J3925-02
Misc :
ALS Vial : 4 Sample Multiplier: 1

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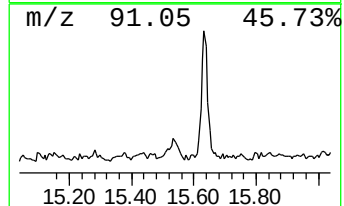
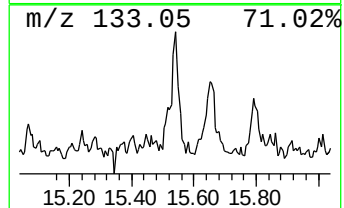
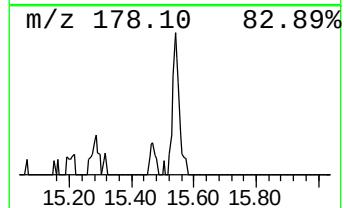
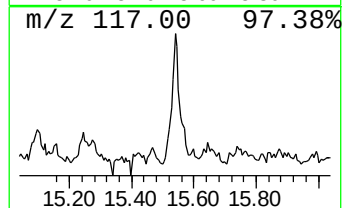
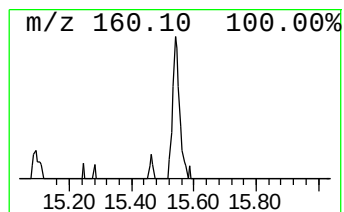
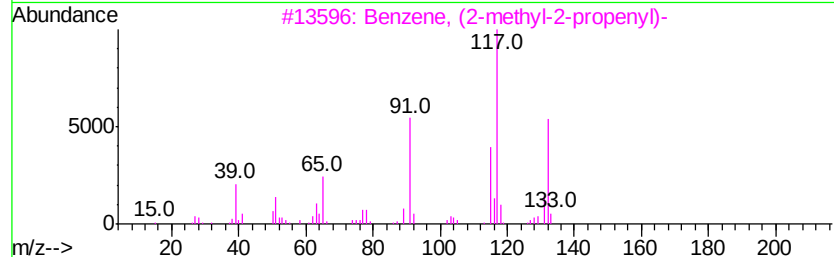
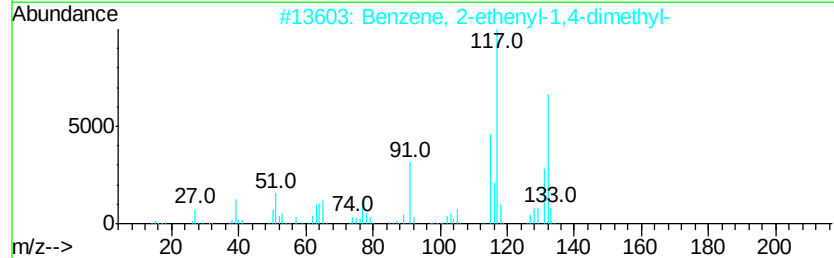
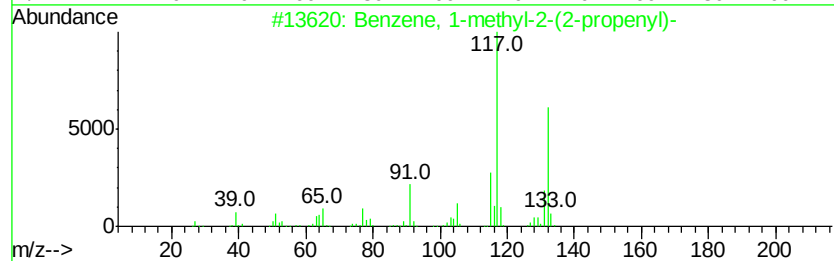
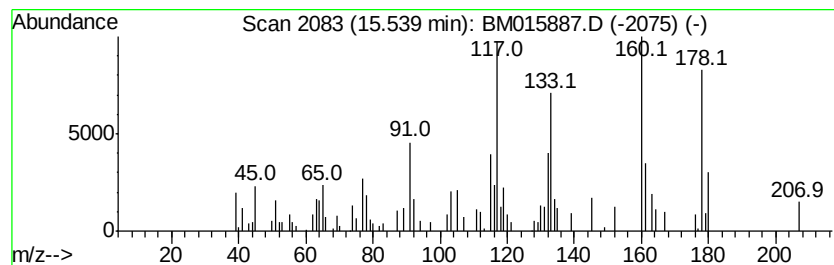
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 unknown15.54 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.54	2.31 ng	113099	Acenaphthene-d10	14.92

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	35
2		Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	35
3		Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	35
4		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	35
5		2,4-Dimethylstyrene	132	C10H12	002234-20-0	35



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM071118\
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Acq On : 11 Jul 2018 14:01
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Sample : J3925-02
Misc :
ALS Vial : 4 Sample Multiplier: 1

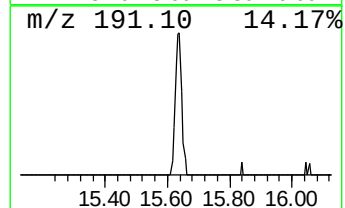
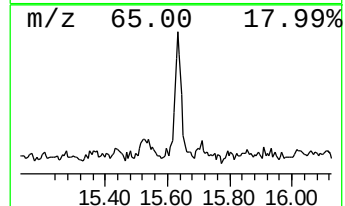
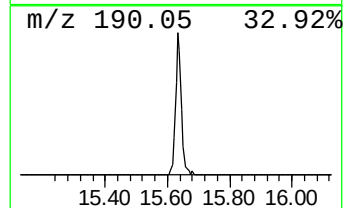
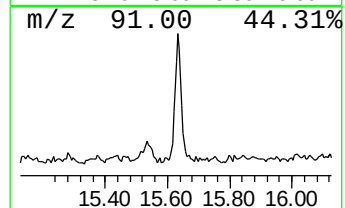
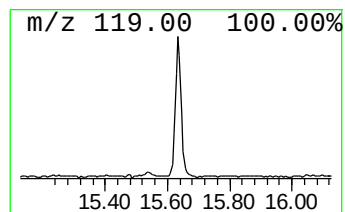
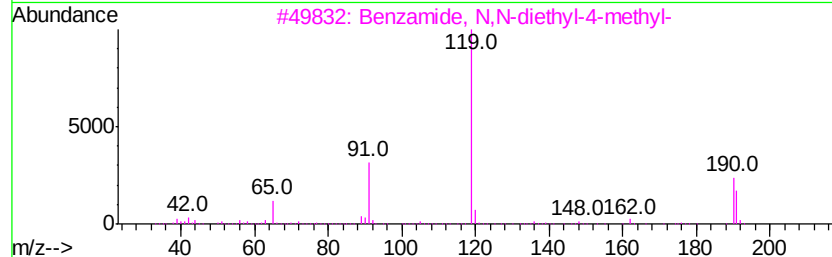
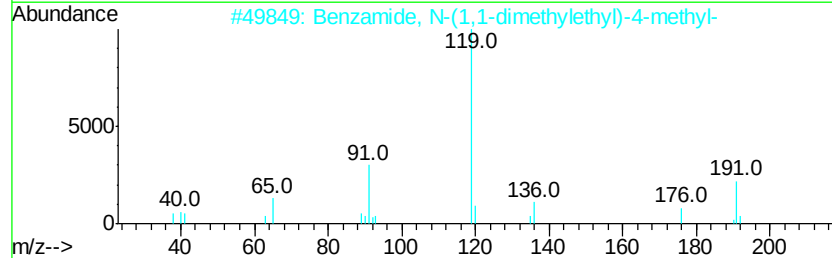
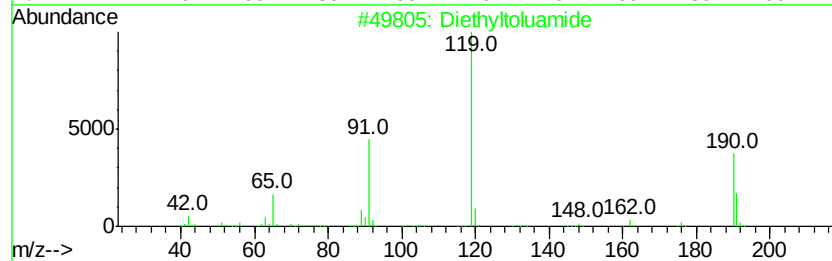
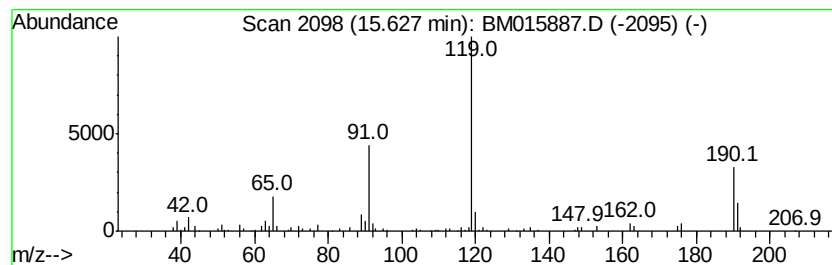
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Diethyltoluamide Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.63	3.25 ng	158842	Acenaphthene-d10	14.92
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Diethyltoluamide	191 C12H17NO	000134-62-3 97
2		Benzamide, N-(1,1-dimethylethyl)...	191 C12H17NO	042498-32-8 83
3		Benzamide, N,N-diethyl-4-methyl-	191 C12H17NO	002728-05-4 80
4		1-Propanone, 2-methyl-1-(4-methy...	162 C11H14O	050390-51-7 64
5		Benzoxazol, 2,3-dihydro-2-thioxo...	269 C15H11NO2S	037442-06-1 59



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM071118\
Data File : BM015887.D
Acq On : 11 Jul 2018 14:01
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Misc :
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ClientSampleId :
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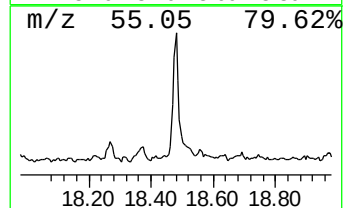
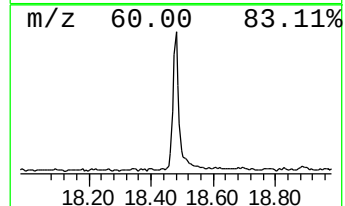
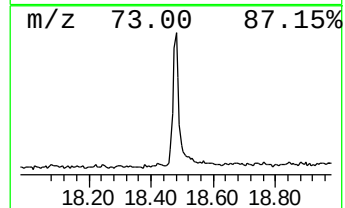
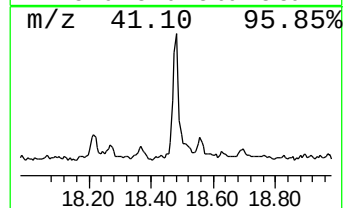
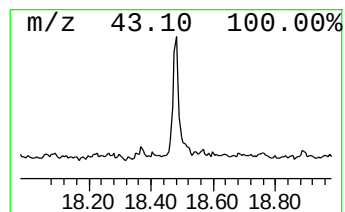
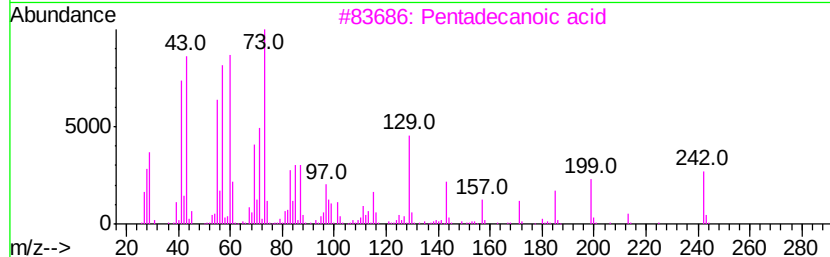
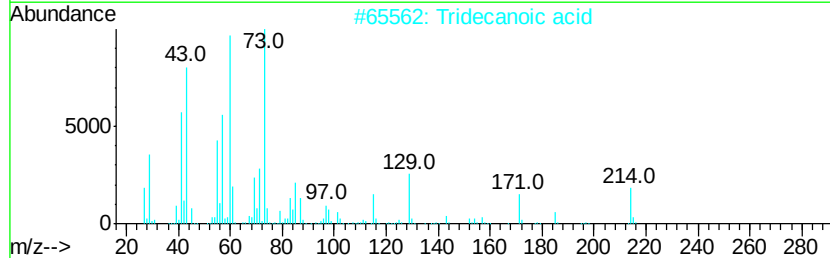
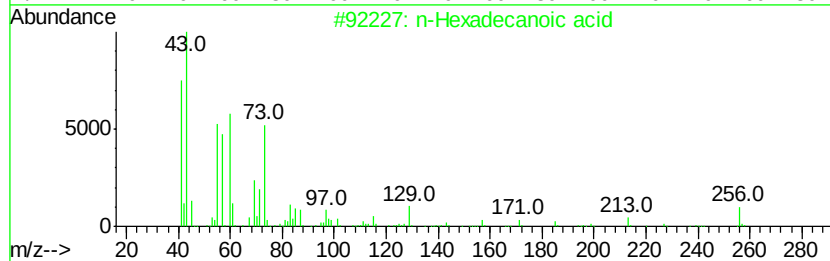
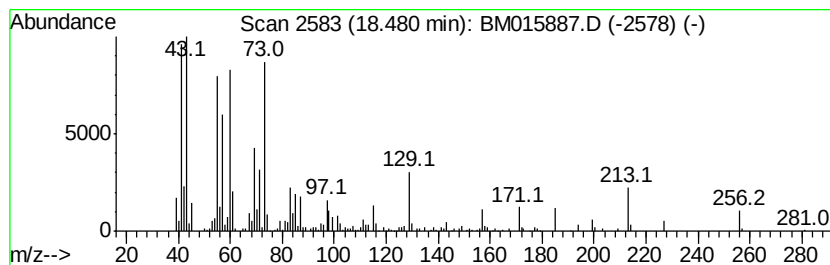
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 n-Hexadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.48	7.14 ng	400706	Phenanthrene-d10	17.64

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
2			Tridecanoic acid	214	C13H26O2	000638-53-9	68
3			Pentadecanoic acid	242	C15H30O2	001002-84-2	68
4			Tetradecanoic acid	228	C14H28O2	000544-63-8	62
5			Dodecanoic acid	200	C12H24O2	000143-07-7	53



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM071118\
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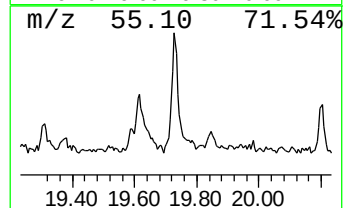
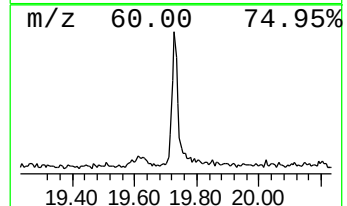
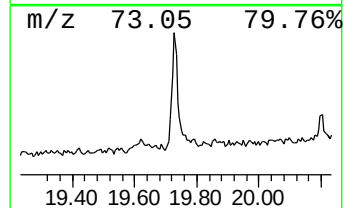
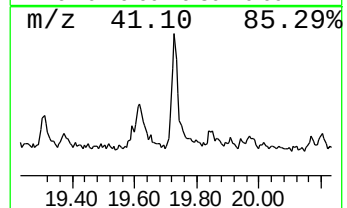
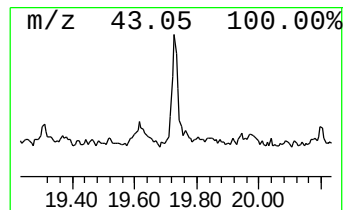
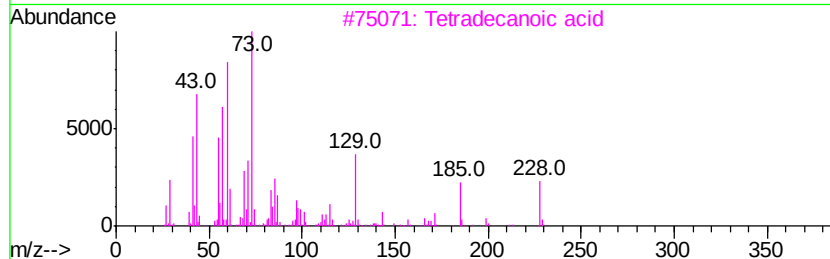
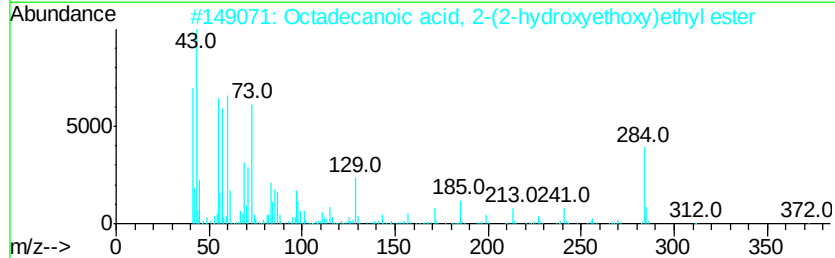
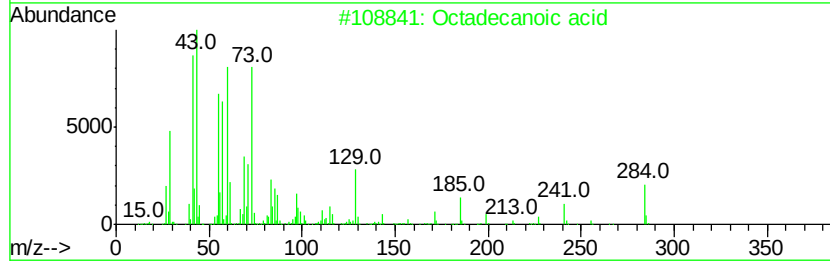
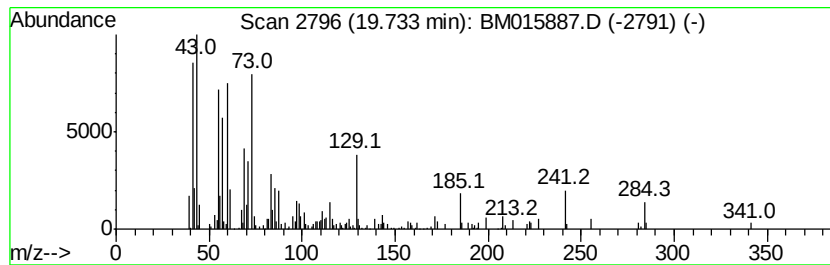
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM070518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Octadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.73	3.81 ng	224157	Chrysene-d12	21.75

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	98
2			Octadecanoic acid, 2-(2-hydroxyve...	372	C22H44O4	000106-11-6	83
3			Tetradecanoic acid	228	C14H28O2	000544-63-8	74
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	64
5			n-Decanoic acid	172	C10H20O2	000334-48-5	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM071118\
Data File : BM015887.D
Acq On : 11 Jul 2018 14:01
Operator : SJ/JU
Sample : J3925-02
Misc :
ALS Vial : 4 Sample Multiplier: 1

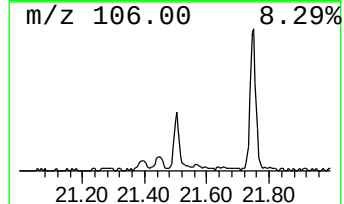
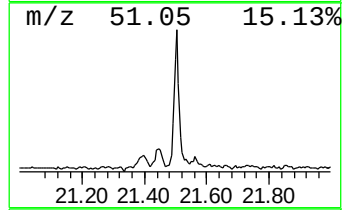
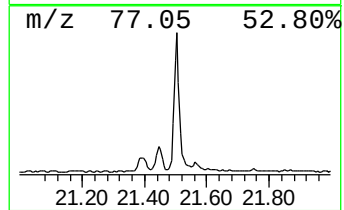
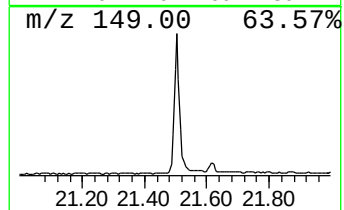
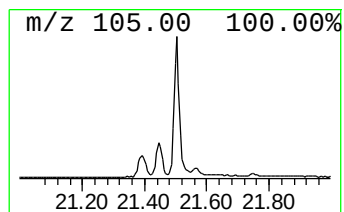
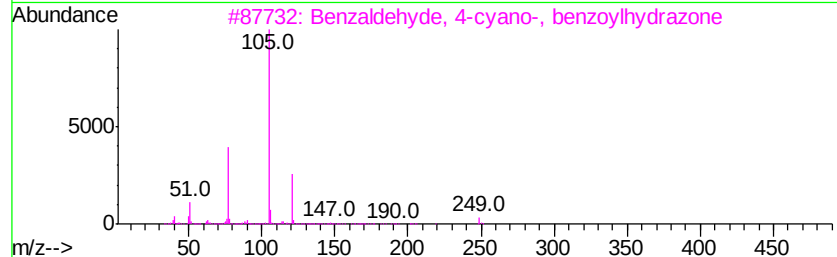
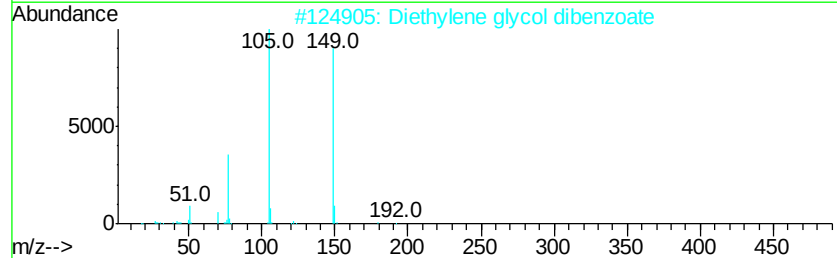
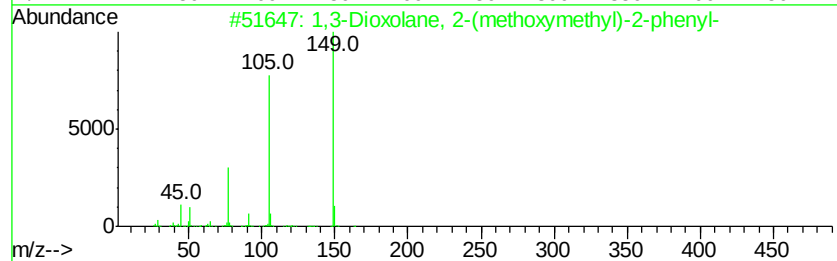
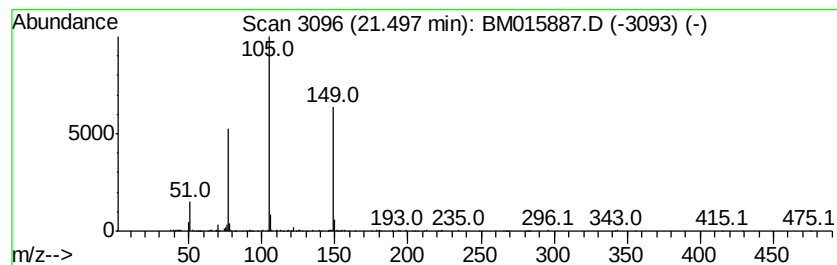
Instrument :
BNA_M
ClientSampleId :
MW-2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM070518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 1,3-Dioxolane, 2-(methoxyme... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.50	6.48 ng	381143	Chrysene-d12	21.75	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Dioxolane, 2-(methoxymethyl)...	194	C11H14O3	1000156-71-2	78
2	Diethylene glycol dibenzoate	314	C18H18O5	000120-55-8	64
3	Benzaldehyde, 4-cyano-, benzoylh...	249	C15H11N3O	195066-12-7	47
4	4-(4-Nitrobenzylidene)-2-phenyl-...	294	C16H10N2O4	007152-75-2	47
5	Benzenebutanoic acid, .gamma.-oxo-	178	C10H10O3	002051-95-8	47



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM071118\
 Data File : BM015887.D
 Acq On : 11 Jul 2018 14:01
 Operator : SJ/JU
 Sample : J3925-02
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 MW-2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\8270-BM070518.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-Butanol, 2,3-di...	3.82	9.1 ng		279770	1	8.29	615061	20.0
2-Pentanone, 4-hy...	5.32	4.2 ng		129824	1	8.29	615061	20.0
unknown7.82	7.82	85.2 ng		2621010	1	8.29	615061	20.0
Ethyl 3,3-dimethy...	8.52	2.4 ng		72551	1	8.29	615061	20.0
unknown9.33	9.33	2.4 ng		74594	1	8.29	615061	20.0
1H-1,2,4-Triazole...	10.12	3.1 ng		128907	2	11.12	830861	20.0
1-Adamantanol	12.36	2.7 ng		110335	2	11.12	830861	20.0
Phenol, 4-(1,1-di...	13.75	33.4 ng		1635150	3	14.92	978796	20.0
unknown15.54	15.54	2.3 ng		113099	3	14.92	978796	20.0
Diethyltoluamide	15.63	3.3 ng		158842	3	14.92	978796	20.0
n-Hexadecanoic acid	18.48	7.1 ng		400706	4	17.64	1122670	20.0
Octadecanoic acid	19.73	3.8 ng		224157	5	21.75	1177160	20.0
1,3-Dioxolane, 2-...	21.50	6.5 ng		381143	5	21.75	1177160	20.0