

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG102518\
 Data File : BG037564.D
 Acq On : 26 Oct 2018 4:07
 Operator : JU/SJ
 Sample : J5633-02
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MW-3

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_G\METHODS\8270-BG101818.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	5.645	394	401	409	rBV	410685	707614	8.54%	2.294%
2	7.243	664	673	688	rBV2	286836	612266	7.39%	1.985%
3	7.625	730	738	750	rBV	798038	1454907	17.55%	4.717%
4	8.101	811	819	831	rBV	201913	372600	4.50%	1.208%
5	8.424	866	874	882	rBV	695190	1269838	15.32%	4.117%
6	9.282	1011	1020	1029	rBV	647893	1204867	14.54%	3.907%
7	10.921	1291	1299	1309	rVB	285132	548292	6.62%	1.778%
8	11.397	1373	1380	1390	rBV	88769	167579	2.02%	0.543%
9	12.737	1597	1608	1612	rBV	3877816	8288143	100.00%	26.874%
10	13.131	1669	1675	1686	rVB	763262	1307777	15.78%	4.240%
11	13.360	1706	1714	1724	rVB	1744882	2794389	33.72%	9.061%
12	14.735	1938	1948	1960	rVB2	491509	859926	10.38%	2.788%
13	14.928	1977	1981	1991	rVB	103657	167427	2.02%	0.543%
14	16.221	2194	2201	2210	rBV2	1362062	2142113	25.85%	6.946%
15	17.120	2348	2354	2359	rBV5	50115	89832	1.08%	0.291%
16	17.484	2407	2416	2422	rBV2	617792	1010271	12.19%	3.276%
17	18.125	2519	2525	2534	rBV3	120455	227831	2.75%	0.739%
18	18.342	2554	2562	2574	rVV2	230820	440586	5.32%	1.429%
19	19.605	2773	2777	2785	rBV2	151523	233403	2.82%	0.757%
20	20.081	2849	2858	2870	rVB	2833531	3865719	46.64%	12.534%
21	21.374	3074	3078	3090	rVB4	69579	160143	1.93%	0.519%
22	21.768	3138	3145	3154	rVB	572593	990940	11.96%	3.213%
23	22.543	3273	3277	3288	rVB3	51230	92744	1.12%	0.301%
24	23.260	3392	3399	3404	rBV	74280	145032	1.75%	0.470%
25	24.088	3534	3540	3549	rVB	81783	191273	2.31%	0.620%
26	25.093	3698	3711	3726	rBV3	331393	1204854	14.54%	3.907%
27	26.233	3898	3905	3916	rVB4	59871	189139	2.28%	0.613%
28	27.637	4138	4144	4155	rVB7	31831	101599	1.23%	0.329%

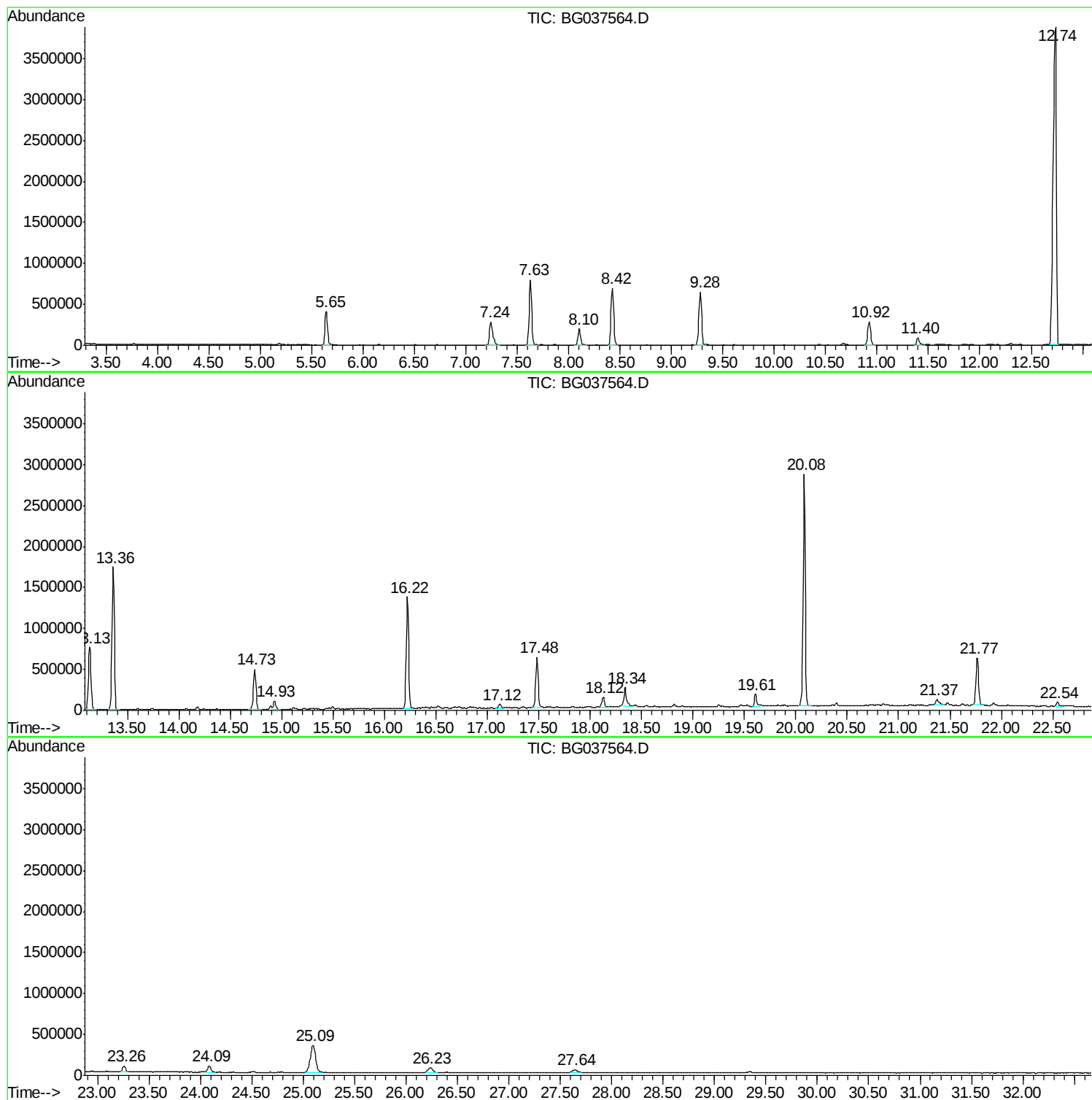
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG101818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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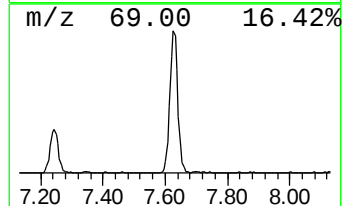
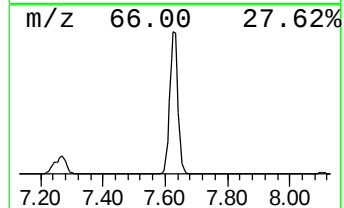
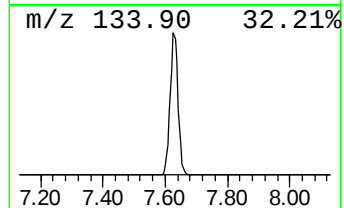
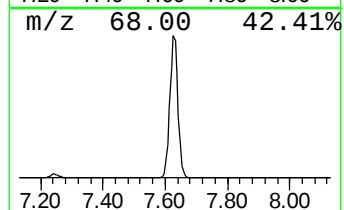
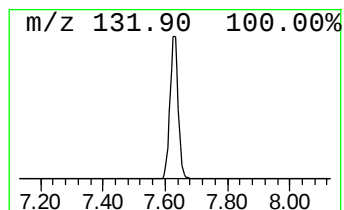
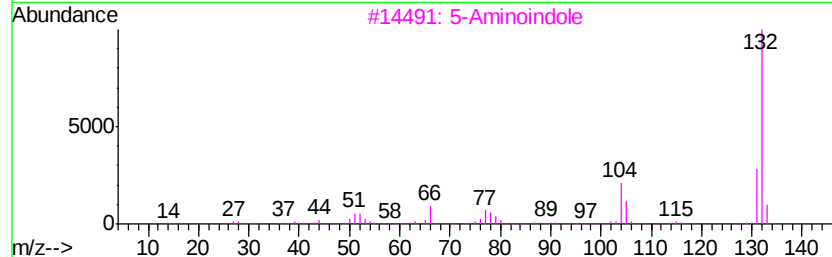
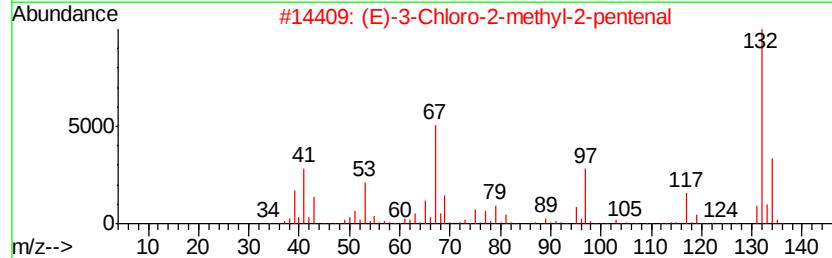
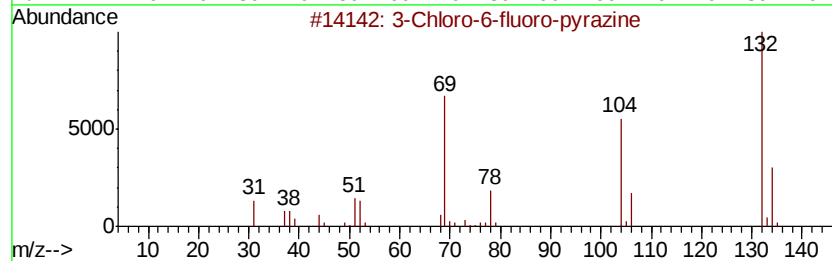
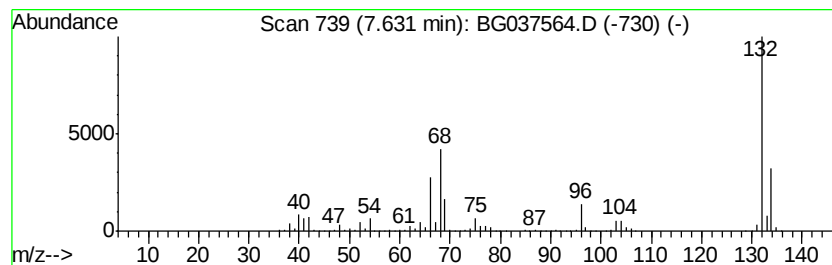
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown7.63 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.63	78.09 ng	1454910	1,4-Dichlorobenzene-d4	8.11

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9



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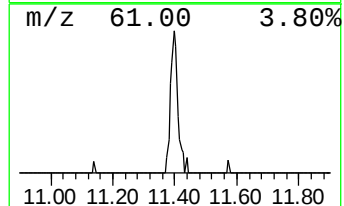
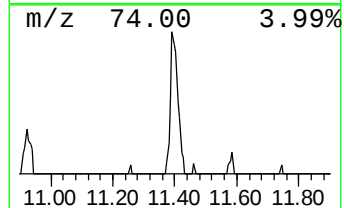
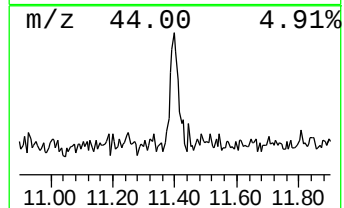
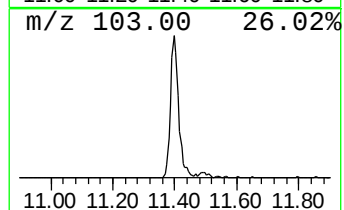
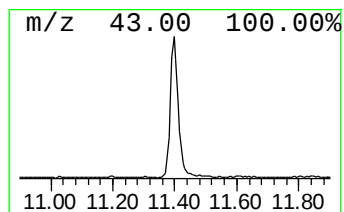
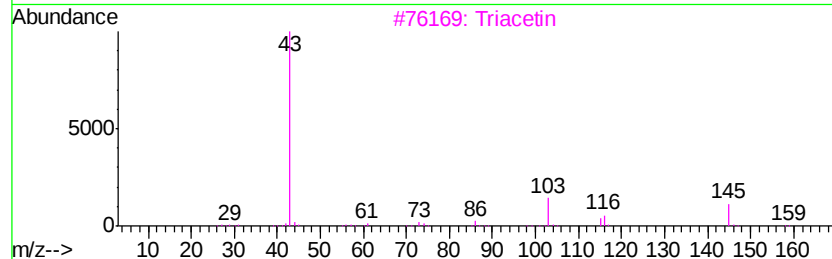
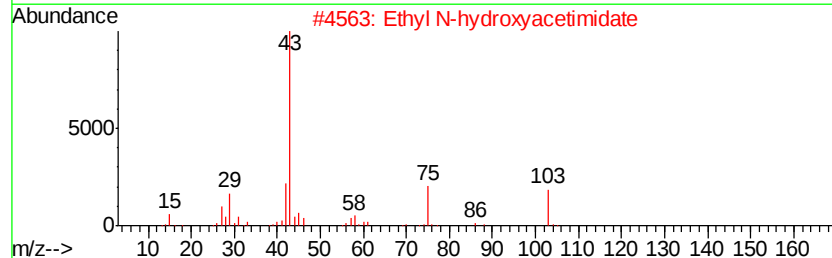
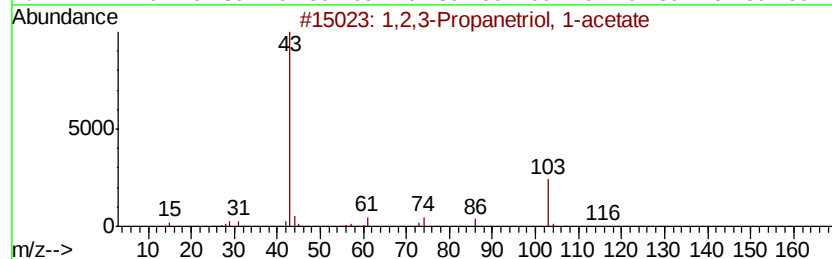
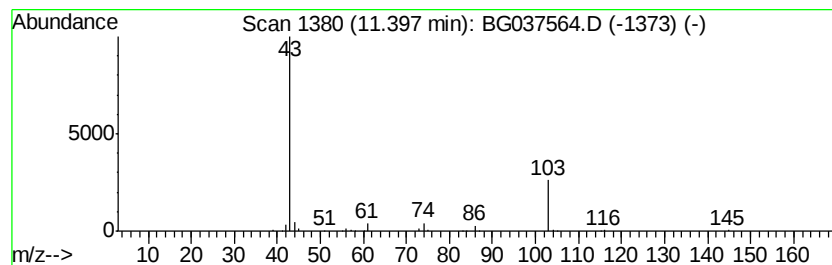
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Peak Number 3 1,2,3-Propanetriol, 1-acetate Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.40	6.11 ng	167579	Naphthalene-d8	10.93

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,2,3-Propanetriol, 1-acetate	134	C5H10O4	000106-61-6	83
2			Ethyl N-hydroxyacetimidate	103	C4H9NO2	010576-12-2	9
3			Triacetin	218	C9H14O6	000102-76-1	9
4			Isobutyl acetate	116	C6H12O2	000110-19-0	9
5			Hydroperoxide, 1-methylbutyl	104	C5H12O2	014018-58-7	5



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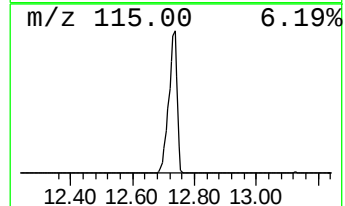
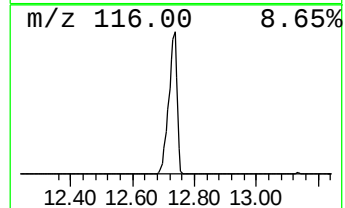
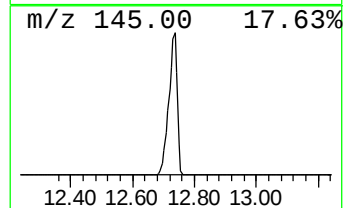
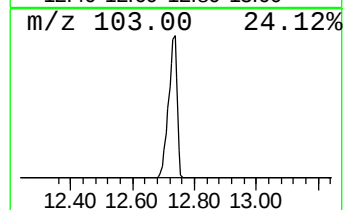
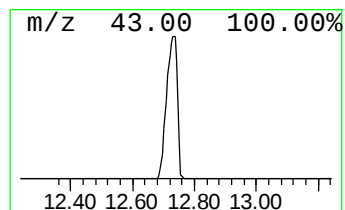
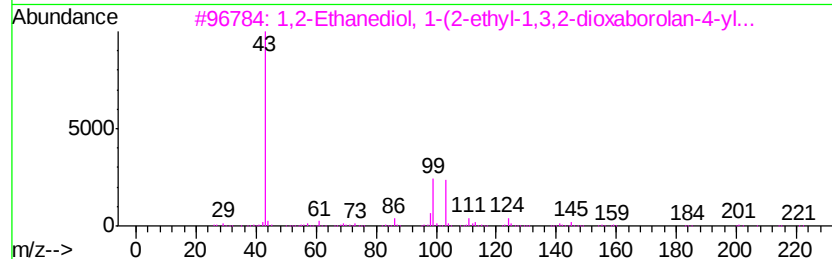
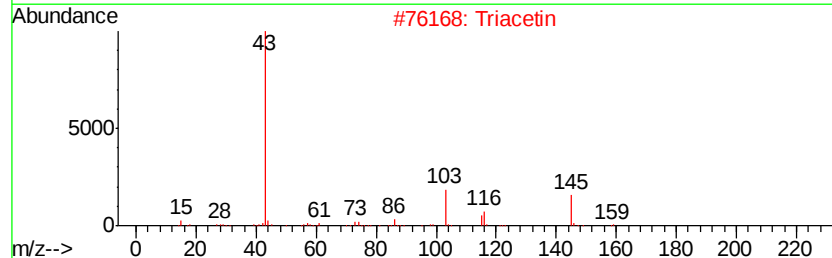
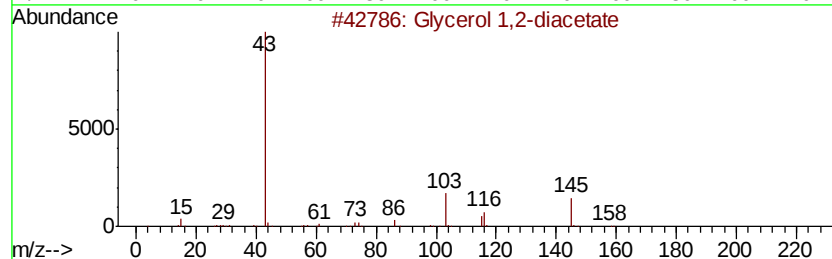
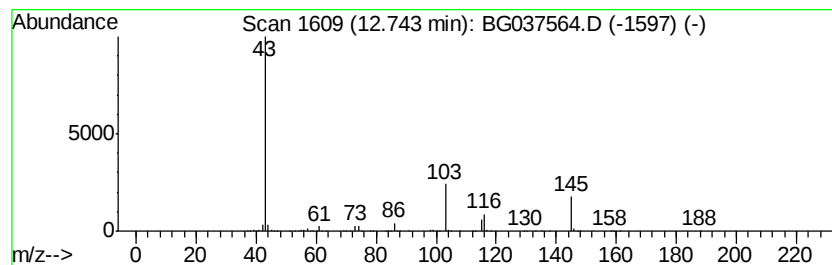
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Peak Number 4 Glycerol 1,2-diacetate Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.74	302.33 ng	8288140	Naphthalene-d8	10.93

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Glycerol 1,2-diacetate	176	C7H12O5	000102-62-5	83
2		Triacetin	218	C9H14O6	000102-76-1	78
3		1,2-Ethenediol, 1-(2-ethyl-1,3,2...	244	C10H17B06	074793-16-1	28
4		1,2,3-Propanetriol, 1-acetate	134	C5H10O4	000106-61-6	25
5		1,2-Epoxy-3-propyl acetate	116	C5H8O3	006387-89-9	10



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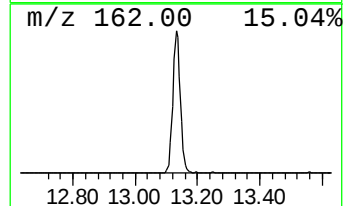
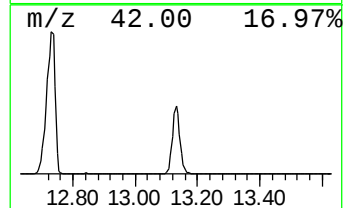
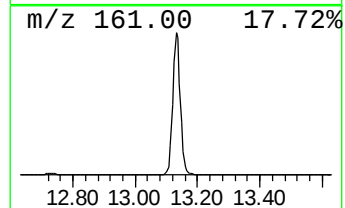
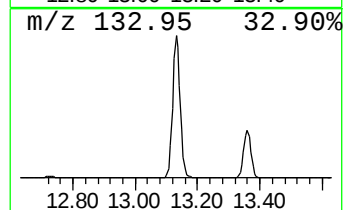
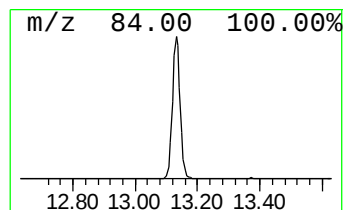
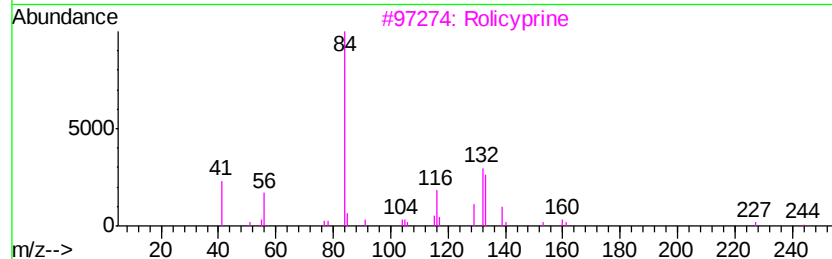
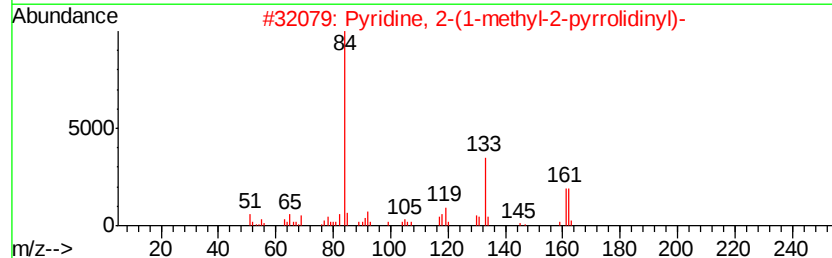
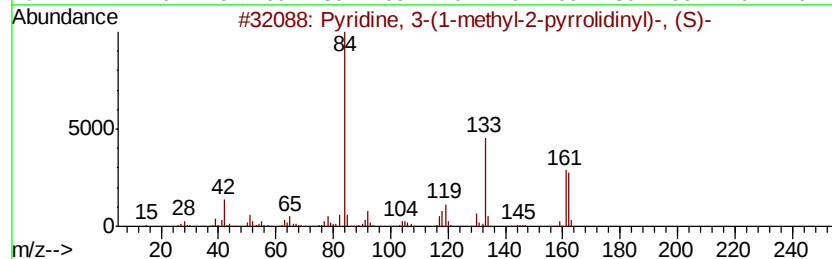
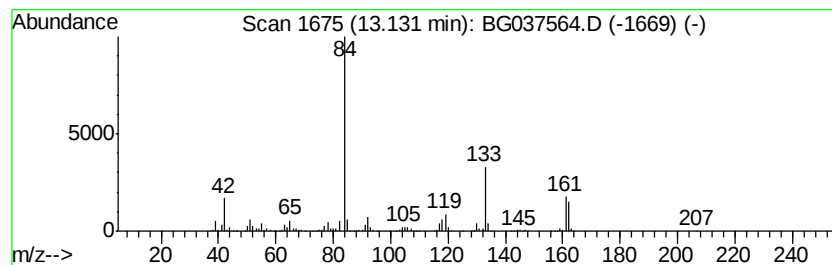
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Peak Number 5 Pyridine, 3-(1-methyl-2-pyr... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.13	30.42 ng	1307780	Acenaphthene-d10	14.73	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyridine, 3-(1-methyl-2-pyrrolid...	162	C10H14N2	000054-11-5	97
2	Pyridine, 2-(1-methyl-2-pyrrolid...	162	C10H14N2	023950-04-1	94
3	Rolicyprine	244	C14H16N2O2	002829-19-8	40
4	2-Methyl[1,3,4]oxadiazole	84	C3H4N2O	003451-51-2	38
5	1-Aminocyclopentane hydroxamic acid	144	C6H12N2O2	062104-33-0	38



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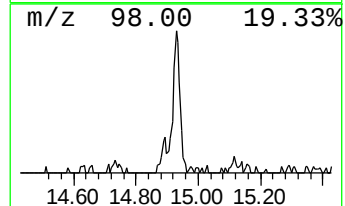
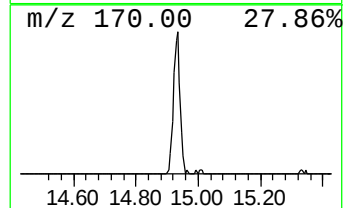
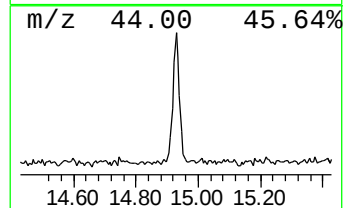
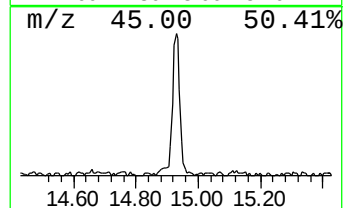
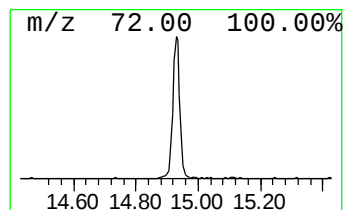
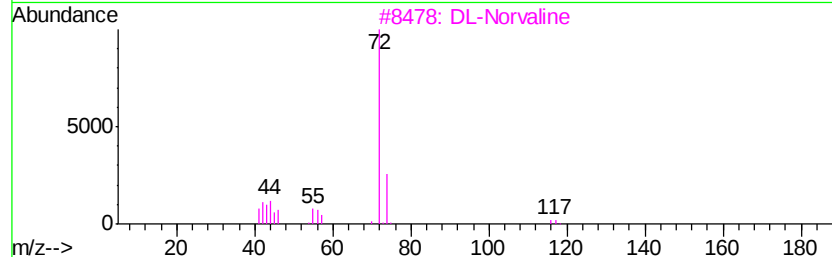
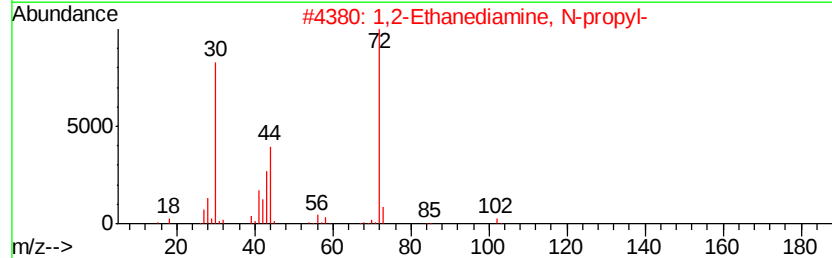
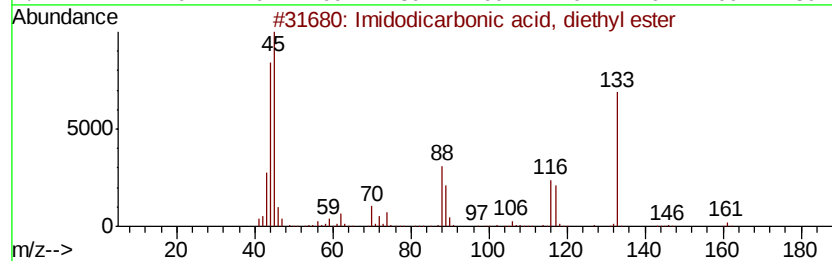
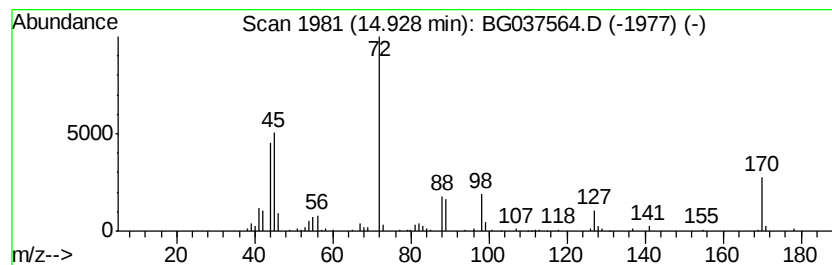
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Peak Number 6 unknown14.93 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.93	3.89 ng	167427	Acenaphthene-d10	14.73

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Imidodicarbonic acid, diethyl ester	161	C6H11N04	019617-44-8	25
2		1,2-Ethanediamine, N-propyl-	102	C5H14N2	000111-39-7	23
3		DL-Norvaline	117	C5H11N02	000760-78-1	12
4		Pentaethylene glycol monododecyl...	406	C22H46O6	003055-95-6	9
5		Butanamide, 3-(2-methoxy-1-oxoet...	307	C15H21N3O4	349575-71-9	9



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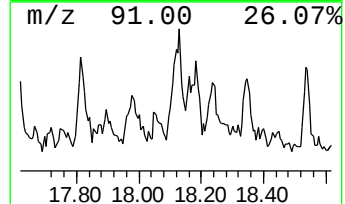
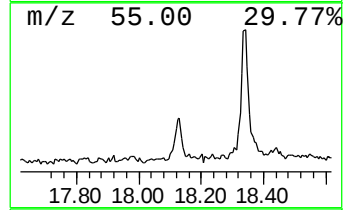
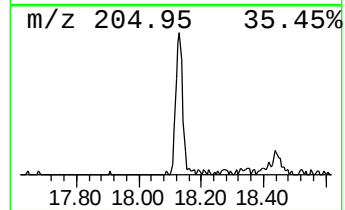
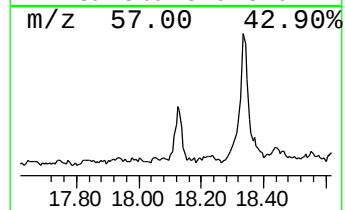
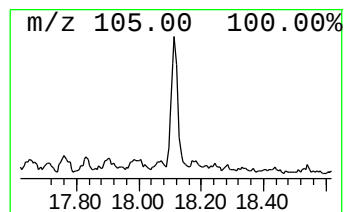
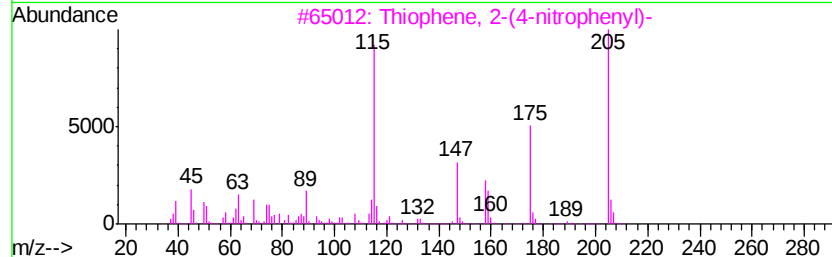
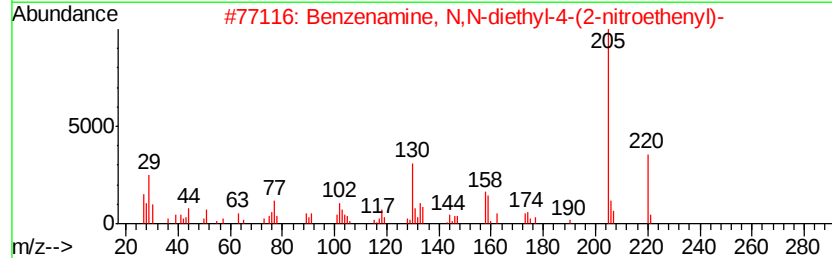
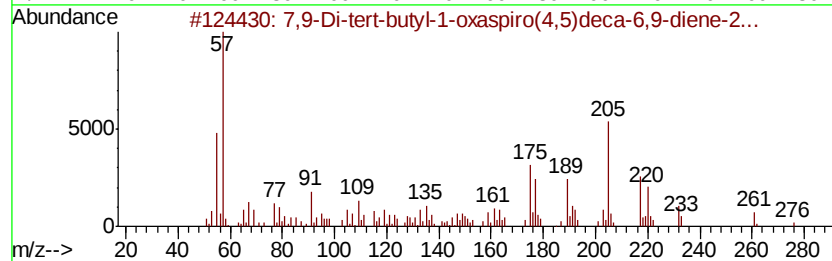
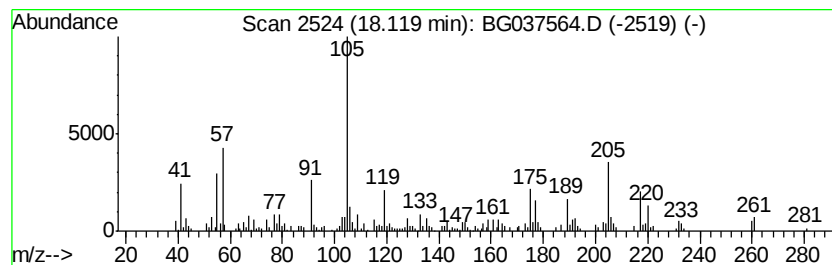
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 7,9-Di-tert-butyl-1-oxaspiro... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.12	4.51 ng	227831	Phenanthrene-d10	17.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	7,9-Di-tert-butyl-1-oxaspiro(4,5...	276	C17H24O3	082304-66-3	92
2		Benzenamine, N,N-diethyl-4-(2-ni...	220	C12H16N2O2	064462-51-7	42
3		Thiophene, 2-(4-nitrophenyl)-	205	C10H7NO2S	059156-21-7	38
4		3-Mercapto-4-phenyl-5-ethyl-1,2,...	205	C10H11N3S	029448-76-8	35
5		2,4-Diamino-6-nitroquinazoline	205	C8H7N5O2	007154-34-9	30



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG102518\
Data File : BG037564.D
Acq On : 26 Oct 2018 4:07
Operator : JU/SJ
Sample : J5633-02
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
MW-3

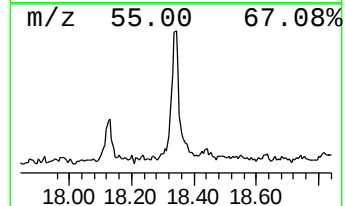
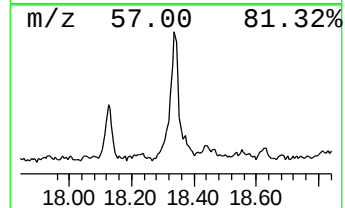
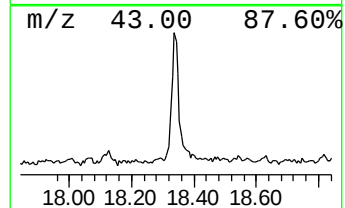
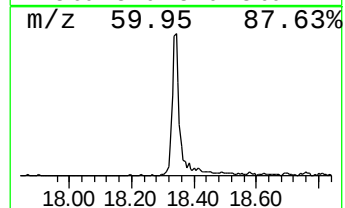
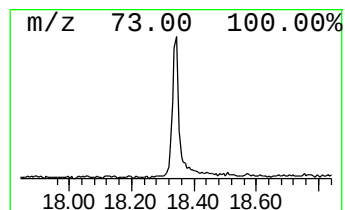
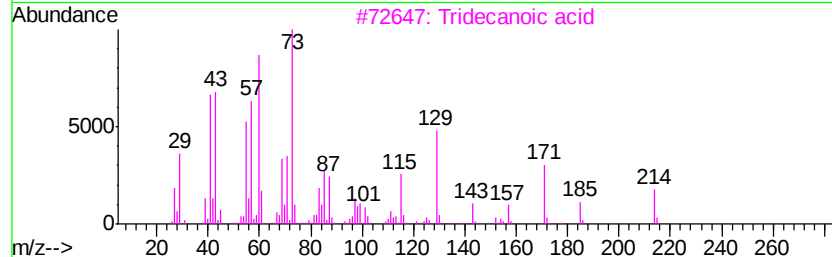
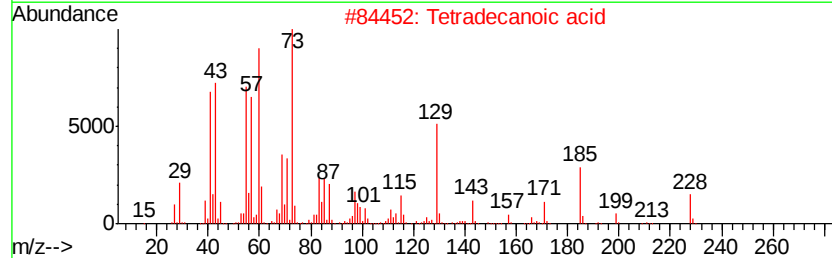
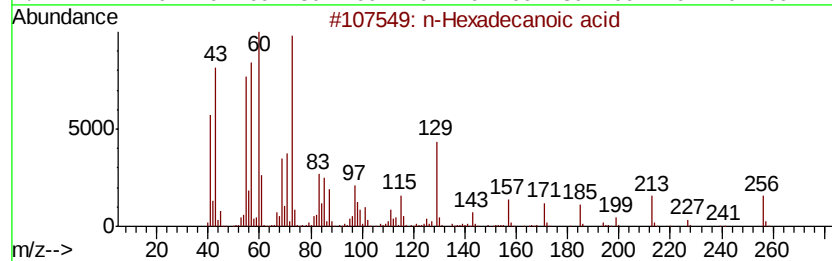
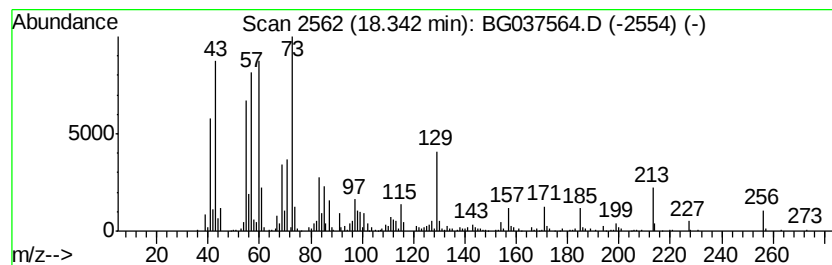
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 n-Hexadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.34	8.72 ng	440586	Phenanthrene-d10	17.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		Tetradecanoic acid	228	C14H28O2	000544-63-8	90
3		Tridecanoic acid	214	C13H26O2	000638-53-9	81
4		Pentadecanoic acid	242	C15H30O2	001002-84-2	81
5		n-Decanoic acid	172	C10H20O2	000334-48-5	78



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG102518\
Data File : BG037564.D
Acq On : 26 Oct 2018 4:07
Operator : JU/SJ
Sample : J5633-02
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
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ClientSampleId :
MW-3

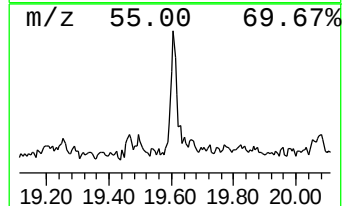
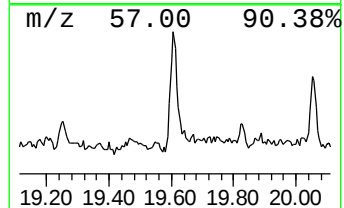
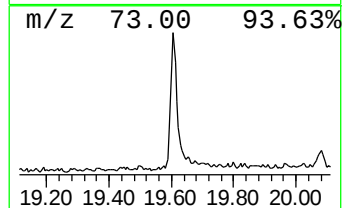
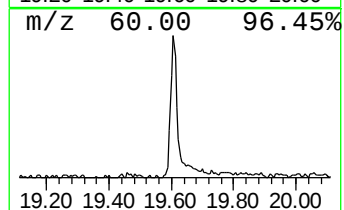
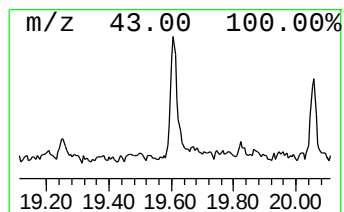
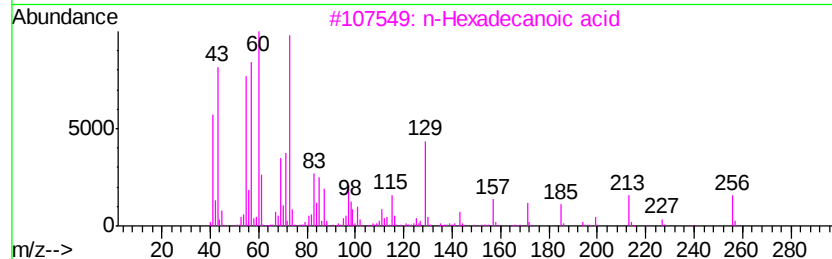
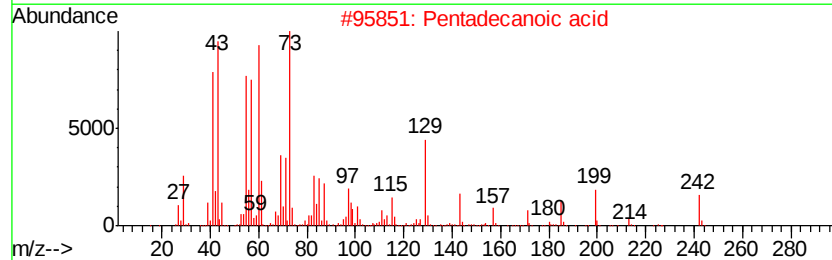
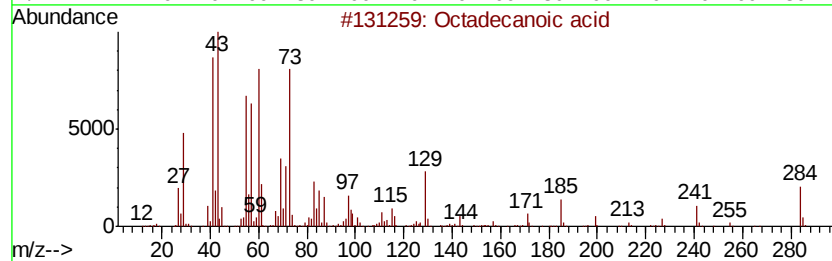
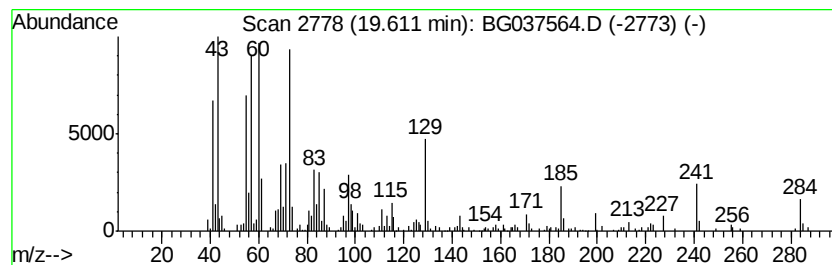
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Octadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.61	4.62 ng	233403	Phenanthrene-d10	17.48

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octadecanoic acid	284	C18H36O2	000057-11-4	98
2		Pentadecanoic acid	242	C15H30O2	001002-84-2	94
3		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	87
4		Tetradecanoic acid	228	C14H28O2	000544-63-8	74
5		n-Decanoic acid	172	C10H20O2	000334-48-5	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG102518\
Data File : BG037564.D
Acq On : 26 Oct 2018 4:07
Operator : JU/SJ
Sample : J5633-02
Misc :
ALS Vial : 20 Sample Multiplier: 1

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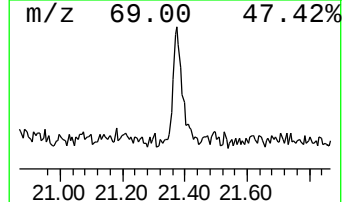
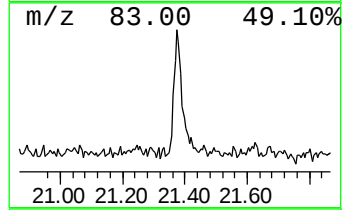
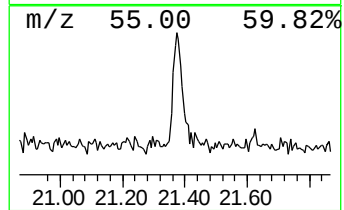
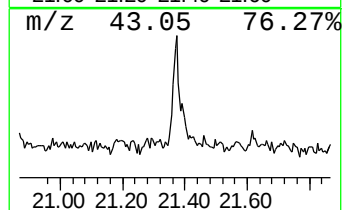
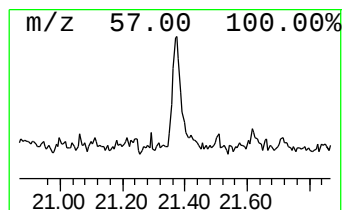
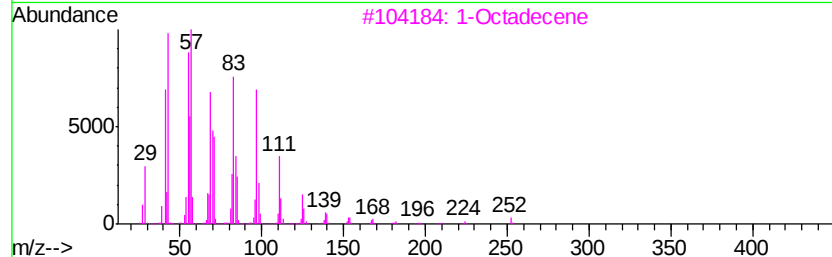
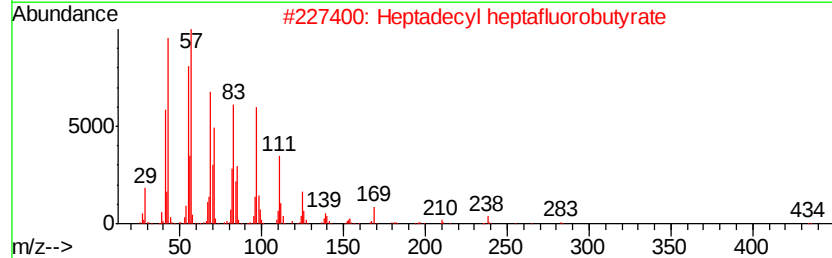
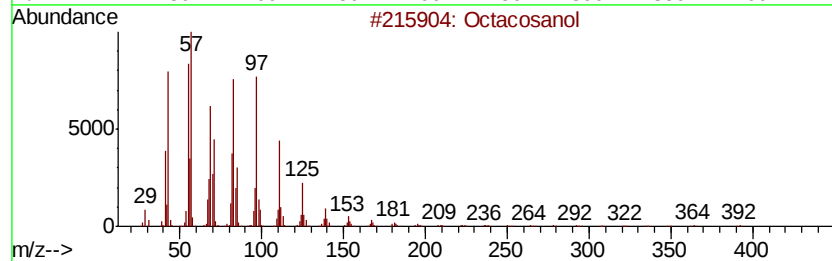
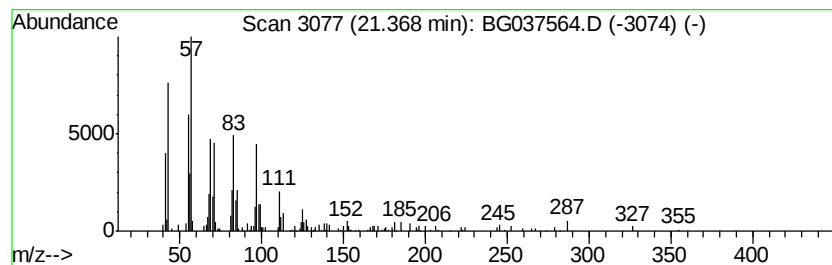
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Octacosanol Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.37	3.23 ng	160143	Chrysene-d12	21.77

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octacosanol	410	C28H58O	000557-61-9	87
2			Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	86
3			1-Octadecene	252	C18H36	000112-88-9	83
4			Triacetyl acetate	480	C32H64O2	041755-58-2	83
5			17-Pentatriacontene	491	C35H70	006971-40-0	83



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG102518\
Data File : BG037564.D
Acq On : 26 Oct 2018 4:07
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Sample : J5633-02
Misc :
ALS Vial : 20 Sample Multiplier: 1

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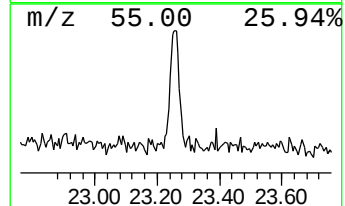
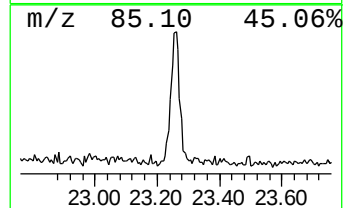
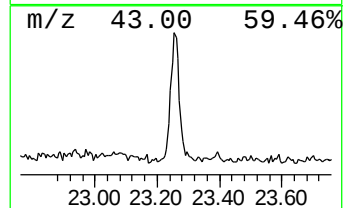
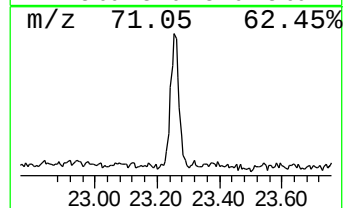
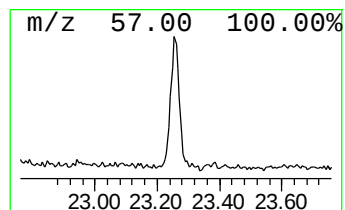
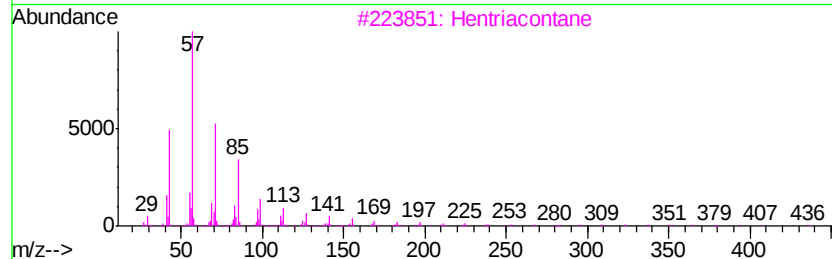
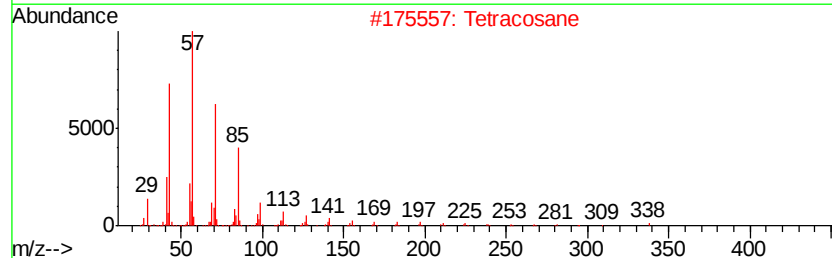
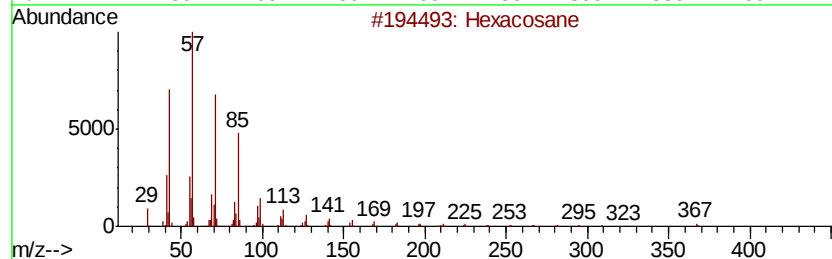
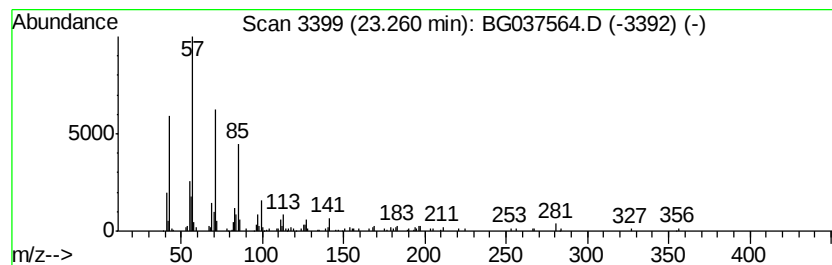
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Hexacosane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.26	2.93 ng	145032	Chrysene-d12	21.77

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexacosane	366	C26H54	000630-01-3	87
2		Tetracosane	338	C24H50	000646-31-1	87
3		Hentriacontane	437	C31H64	000630-04-6	87
4		Octacosane	394	C28H58	000630-02-4	87
5		Heptacosane	380	C27H56	000593-49-7	83



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG102518\
Data File : BG037564.D
Acq On : 26 Oct 2018 4:07
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Sample : J5633-02
Misc :
ALS Vial : 20 Sample Multiplier: 1

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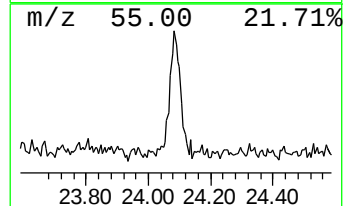
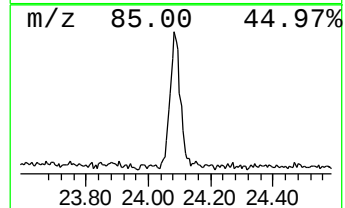
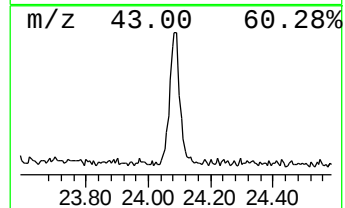
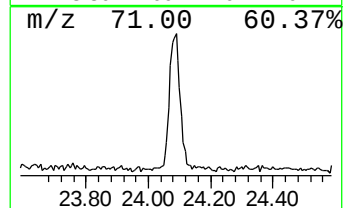
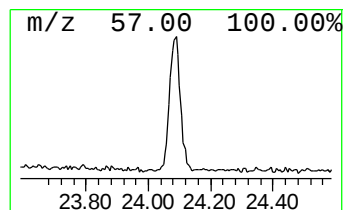
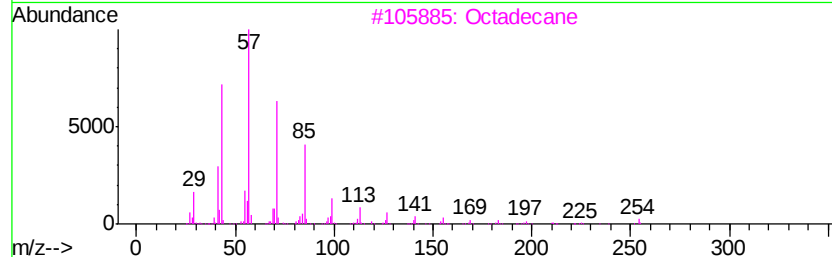
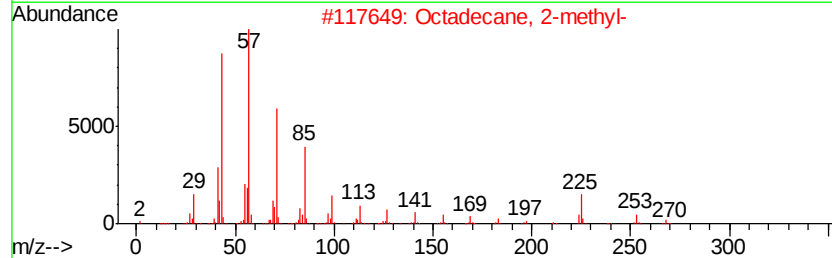
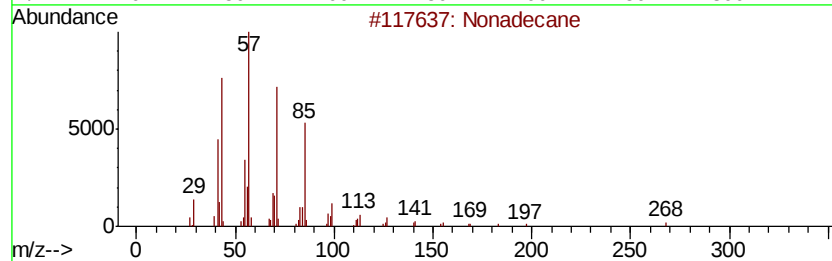
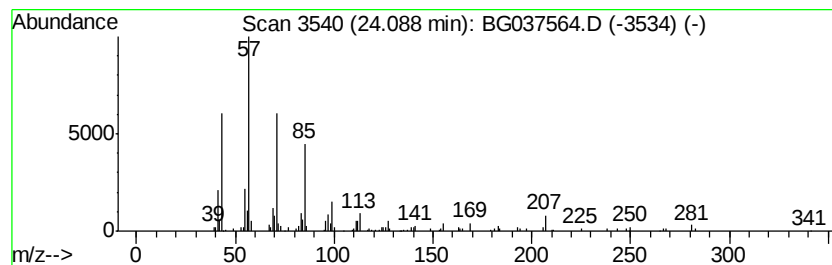
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Nonadecane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.09	3.18 ng	191273	Perylene-d12	25.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonadecane	268	C19H40	000629-92-5	90
2		Octadecane, 2-methyl-	268	C19H40	001560-88-9	87
3		Octadecane	254	C18H38	000593-45-3	87
4		Octacosane	394	C28H58	000630-02-4	87
5		Heptacosane	380	C27H56	000593-49-7	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG102518\
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ClientSampleId :
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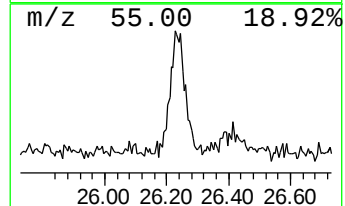
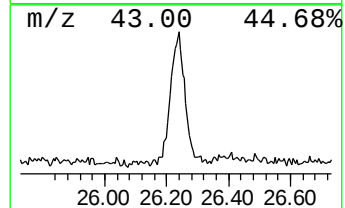
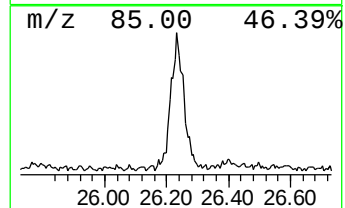
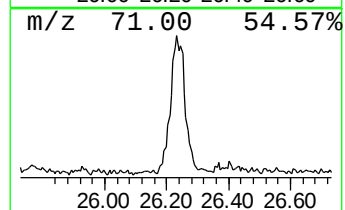
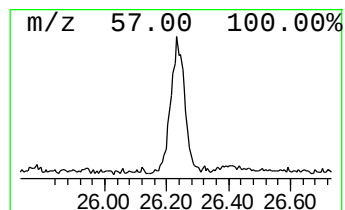
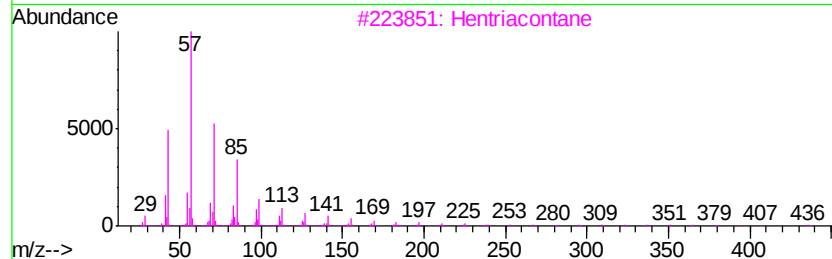
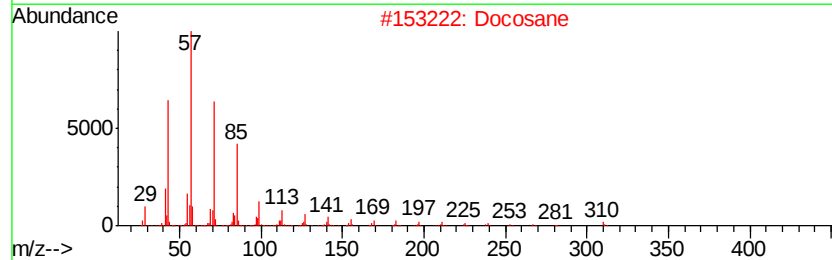
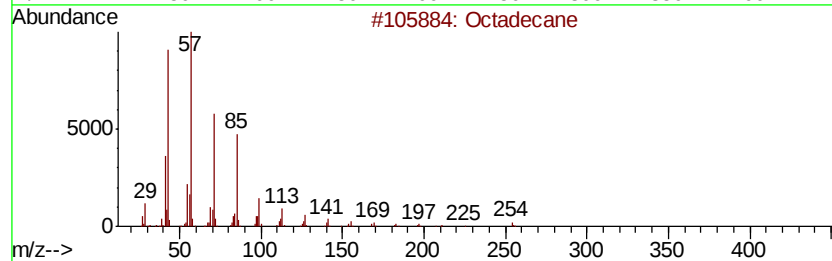
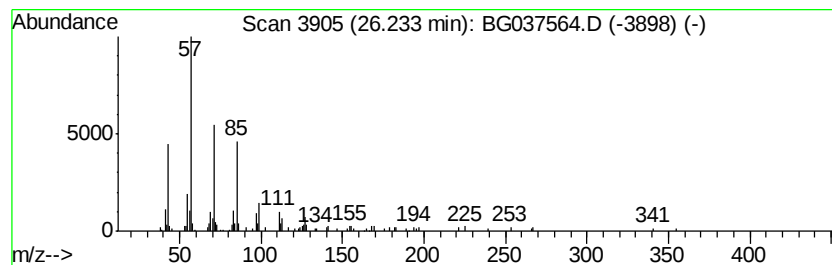
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Octadecane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.23	3.14 ng	189139	Perylene-d12	25.10

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecane	254	C18H38	000593-45-3	72
2			Docosane	310	C22H46	000629-97-0	64
3			Hentriacontane	437	C31H64	000630-04-6	64
4			Heptacosane	380	C27H56	000593-49-7	58
5			Hexacosane	366	C26H54	000630-01-3	52



Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG102518\
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Acq On : 26 Oct 2018 4:07
Operator : JU/SJ
Sample : J5633-02
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG101818.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
unknown7.63	7.63	78.1	ng	1454910	1	8.11	372600	20.0
1,2,3-Propanetriol	11.40	6.1	ng	167579	2	10.93	548292	20.0
Glycerol 1,2-diac...	12.74	302.3	ng	8288140	2	10.93	548292	20.0
Pyridine, 3-(1-me...	13.13	30.4	ng	1307780	3	14.73	859926	20.0
unknown14.93	14.93	3.9	ng	167427	3	14.73	859926	20.0
7,9-Di-tert-butyl...	18.12	4.5	ng	227831	4	17.48	1010270	20.0
n-Hexadecanoic acid	18.34	8.7	ng	440586	4	17.48	1010270	20.0
Octadecanoic acid	19.61	4.6	ng	233403	4	17.48	1010270	20.0
Octacosanol	21.37	3.2	ng	160143	5	21.77	990940	20.0
Hexacosane	23.26	2.9	ng	145032	5	21.77	990940	20.0
Nonadecane	24.09	3.2	ng	191273	6	25.10	1204850	20.0
Octadecane	26.23	3.1	ng	189139	6	25.10	1204850	20.0