

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM060119\
 Data File : BM020650.D
 Acq On : 01 Jun 2019 14:56
 Operator : HP/JU
 Sample : K3075-09
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 TEST-PIT-5

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-BM053119.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	3.064	8	13	30	rVB	2253738	2620431	18.96%	2.540%
2	5.410	405	412	426	rBV	5490311	8160536	59.06%	7.910%
3	6.993	671	681	696	rBV	5641893	9317311	67.43%	9.031%
4	7.363	736	744	754	rBV	5773040	9175680	66.40%	8.894%
5	7.828	816	823	830	rBV	1087713	1711382	12.38%	1.659%
6	8.152	868	878	885	rBV	3985137	6442058	46.62%	6.244%
7	8.993	1012	1021	1031	rBV	3413822	5580582	40.39%	5.409%
8	10.622	1289	1298	1307	rBV	1428855	2355661	17.05%	2.283%
9	11.640	1465	1471	1481	rVB	229701	401113	2.90%	0.389%
10	13.086	1710	1717	1726	rBV	7438203	11000954	79.61%	10.663%
11	13.922	1853	1859	1868	rBV	889043	1249660	9.04%	1.211%
12	14.469	1941	1952	1959	rBV2	2012477	3134261	22.68%	3.038%
13	15.957	2198	2205	2222	rBV	6182369	9243226	66.89%	8.959%
14	17.210	2412	2418	2425	rBV2	2677693	3658667	26.48%	3.546%
15	18.110	2565	2571	2581	rBV	914447	1342987	9.72%	1.302%
16	19.386	2783	2788	2800	rBV	408454	631806	4.57%	0.612%
17	19.839	2858	2865	2870	rBV	11015837	13818433	100.00%	13.394%
18	21.104	3076	3080	3089	rBV2	857260	1124110	8.13%	1.090%
19	21.386	3123	3128	3134	rBV	2965565	3616302	26.17%	3.505%
20	21.551	3152	3156	3161	rVB	271434	321649	2.33%	0.312%
21	21.992	3227	3231	3236	rBV	464394	604512	4.37%	0.586%
22	22.468	3308	3312	3317	rBV	648544	817363	5.92%	0.792%
23	22.604	3331	3335	3340	rVB	252997	337774	2.44%	0.327%
24	22.992	3397	3401	3407	rBV	634620	899945	6.51%	0.872%
25	23.586	3496	3502	3509	rVB	572365	969447	7.02%	0.940%
26	23.680	3511	3518	3527	rVB	1809678	3473915	25.14%	3.367%
27	24.256	3611	3616	3623	rVB	385790	741617	5.37%	0.719%
28	25.039	3745	3749	3758	rVB2	194530	416638	3.02%	0.404%

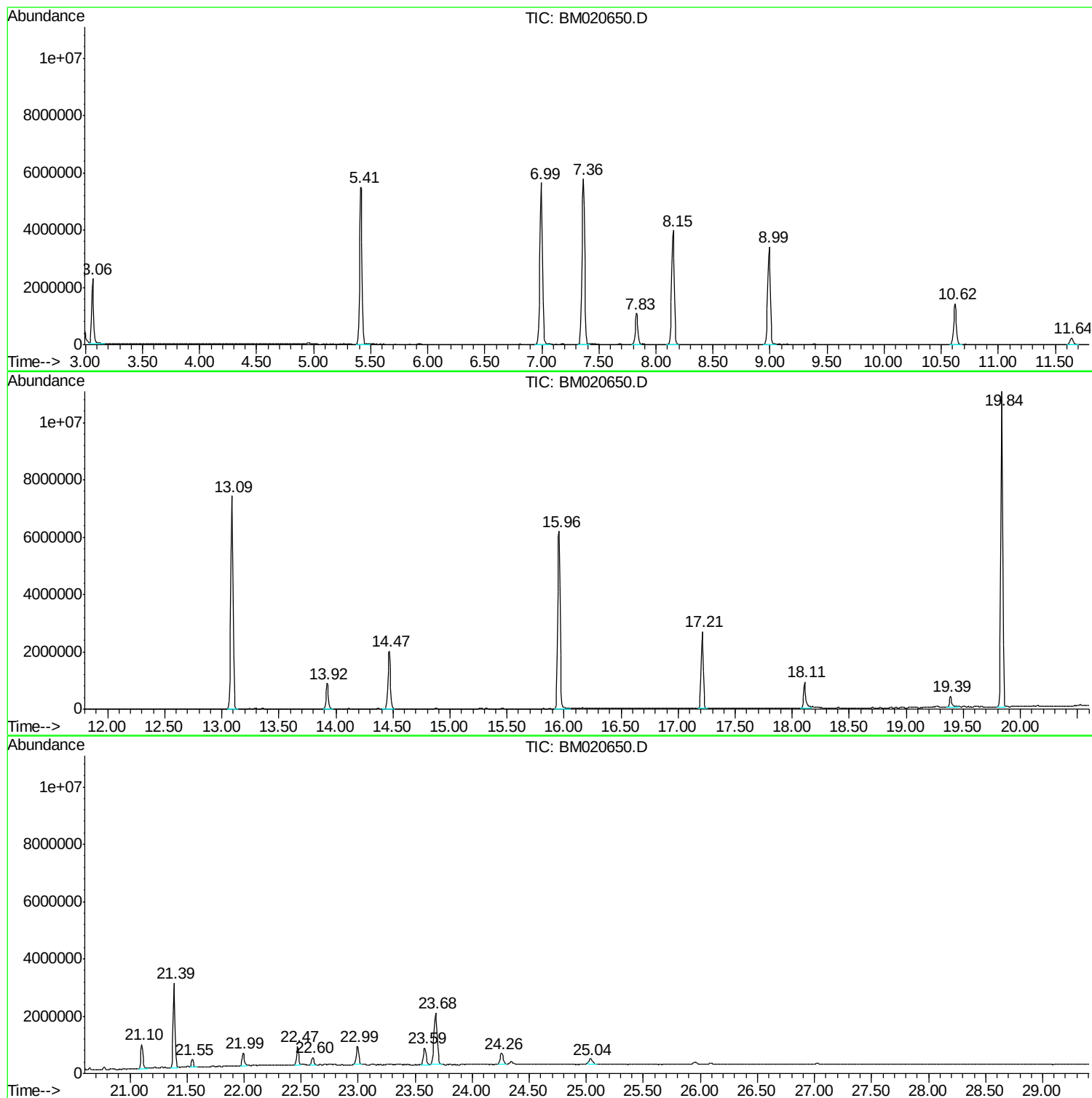
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ClientSampled :
TEST-PIT-5

Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM053119.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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Instrument :
BNA_M
ClientSampleId :
TEST-PIT-5

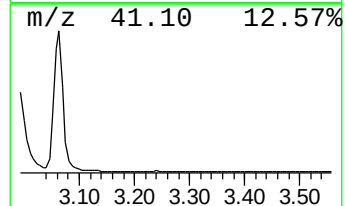
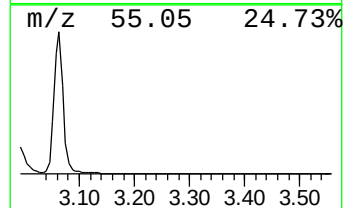
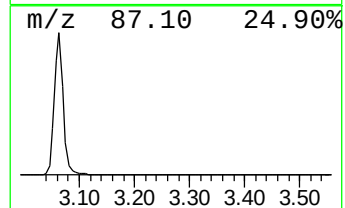
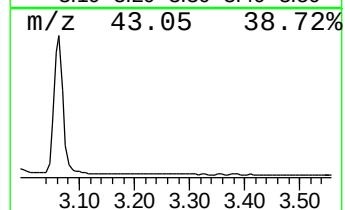
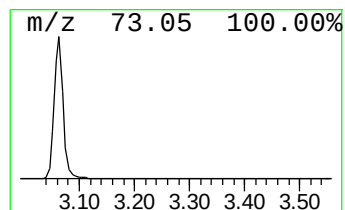
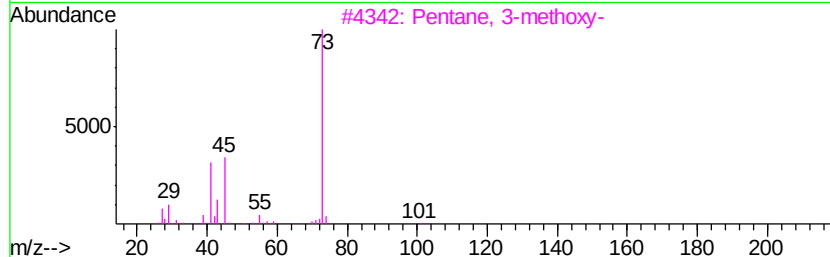
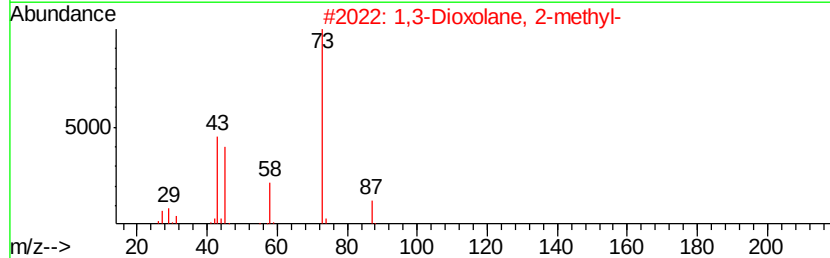
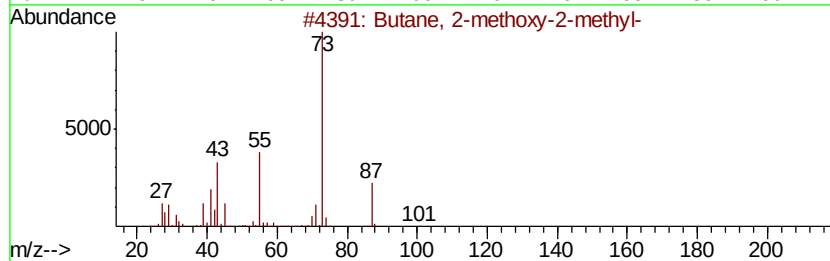
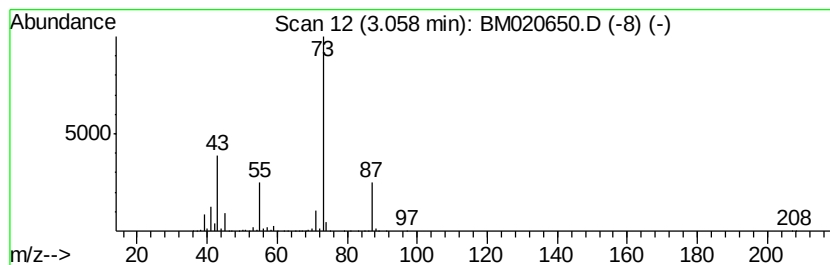
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM053119.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.06	30.62 ng	2620430	1,4-Dichlorobenzene-d4	7.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	37
3		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
4		Silane, tetramethyl-	88	C4H12Si	000075-76-3	17
5		N-Ethylformamide	73	C3H7NO	000627-45-2	9



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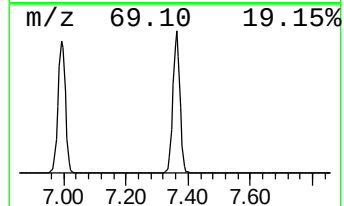
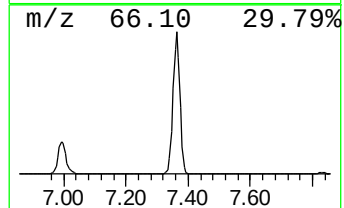
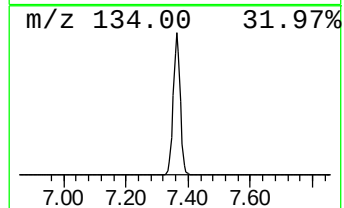
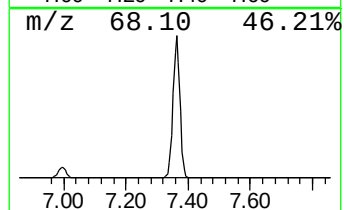
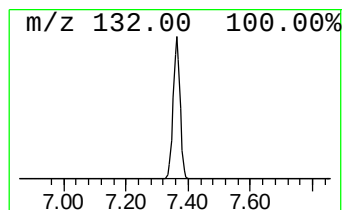
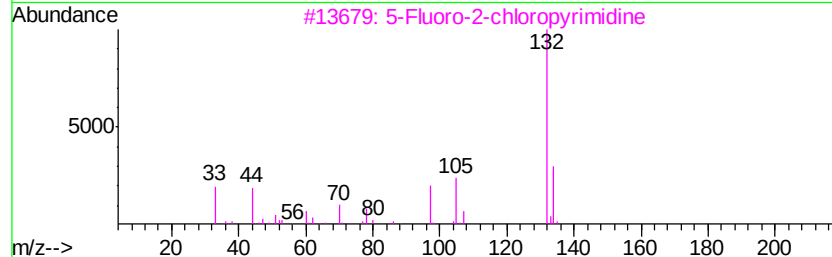
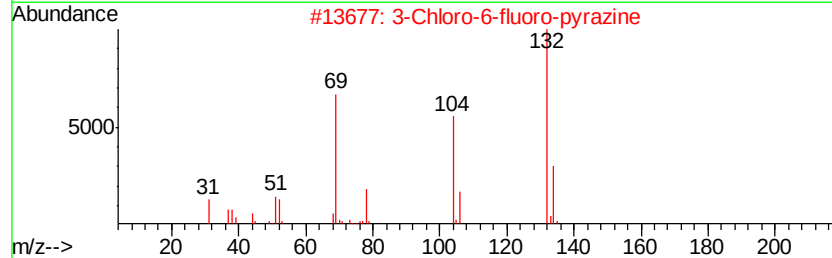
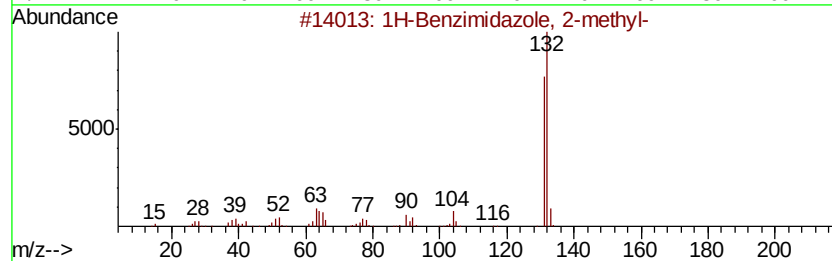
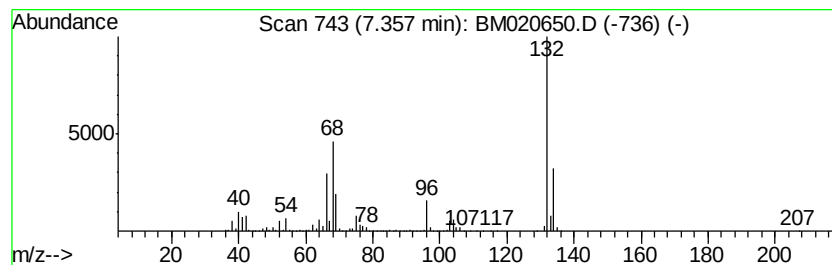
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown7.36 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.36	107.23 ng	9175680	1,4-Dichlorobenzene-d4	7.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
3		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
5		Benzene, 1,2,4-trifluoro-	132	C6H3F3	000367-23-7	9



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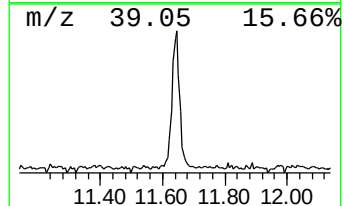
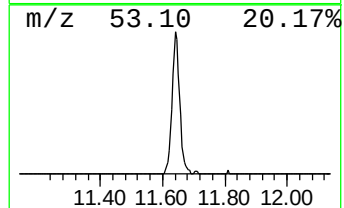
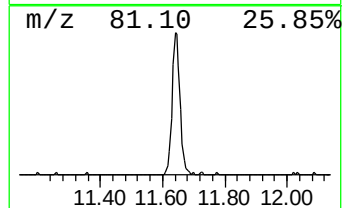
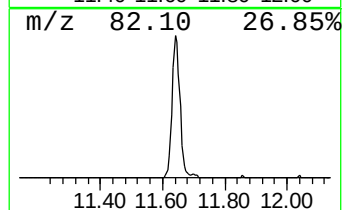
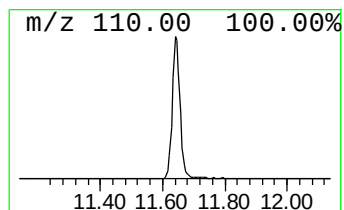
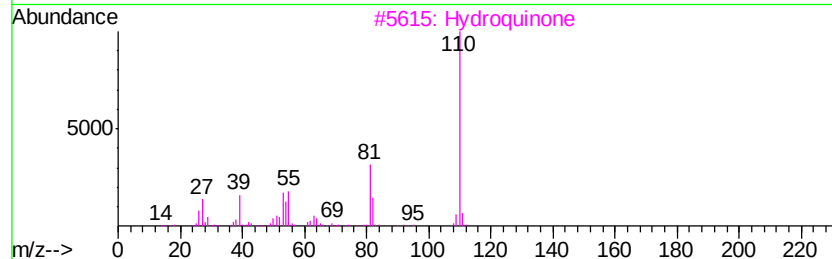
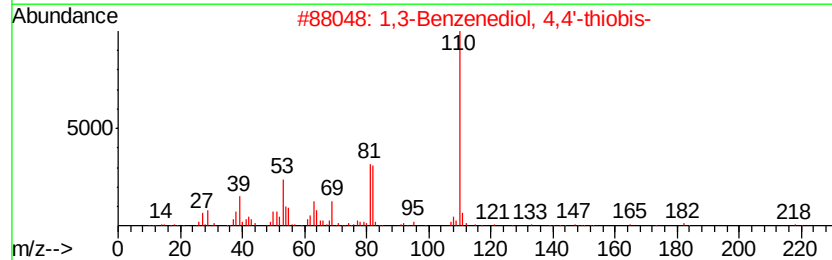
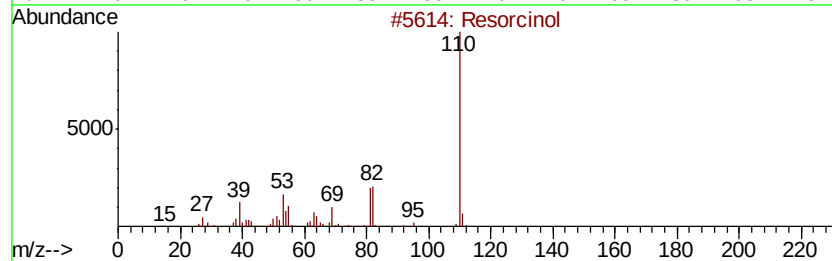
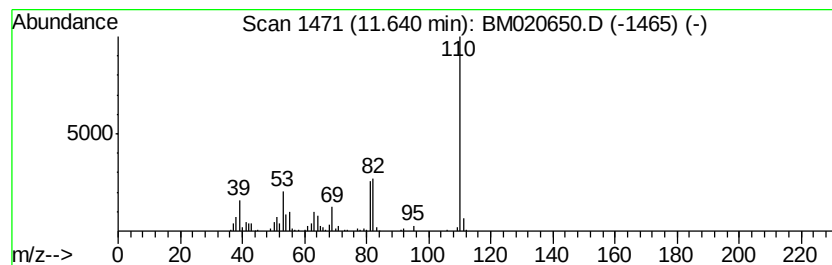
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Peak Number 4 Resorcinol Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.64	3.41 ng	401113	Napthalene-d8	10.62
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Resorcinol		110 C6H6O2	000108-46-3 97
2	1,3-Benzenediol, 4,4'-thiobis-		250 C12H10O4S	000097-29-0 90
3	Hydroquinone		110 C6H6O2	000123-31-9 80
4	3(2H)-Pyridazinone, 6-methyl-		110 C5H6N2O	013327-27-0 53
5	1H-Imidazole-2-carboxaldehyde, 1...		110 C5H6N2O	013750-81-7 52



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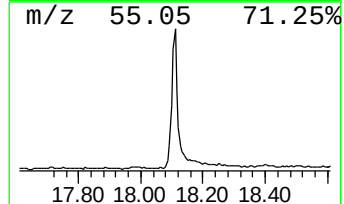
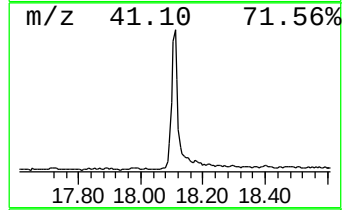
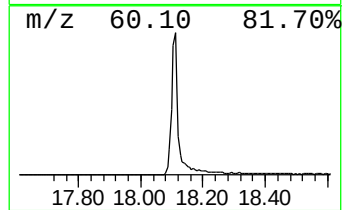
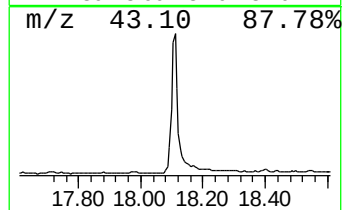
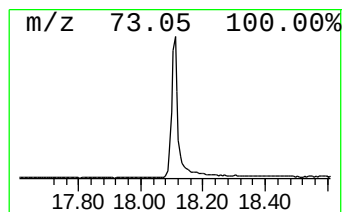
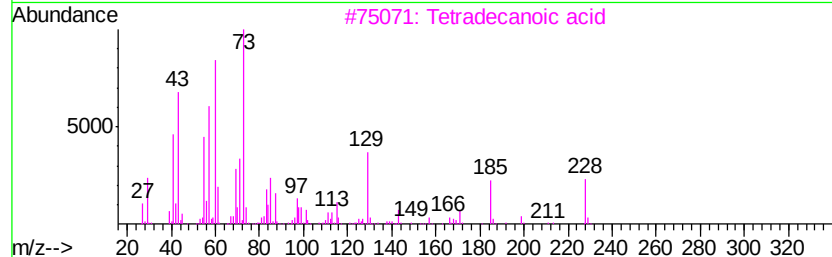
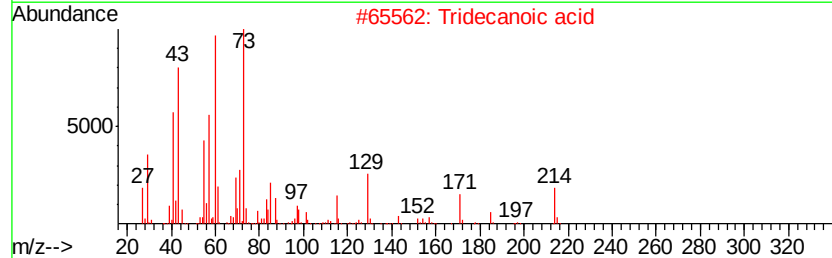
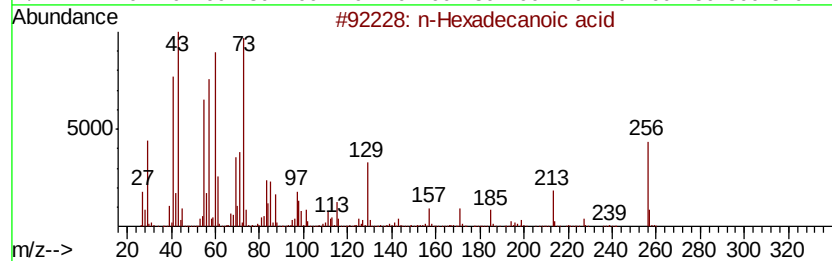
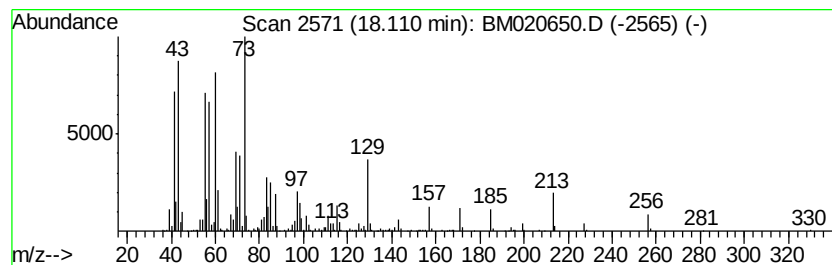
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 5 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD		R.T.
18.11	7.34 ng	1342990	Phenanthrene-d10		17.21
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2	Tridecanoic acid	214	C13H26O2	000638-53-9	91
3	Tetradecanoic acid	228	C14H28O2	000544-63-8	80
4	Undecane	156	C11H24	001120-21-4	11
5	Dodecane	170	C12H26	000112-40-3	11



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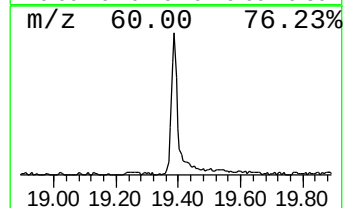
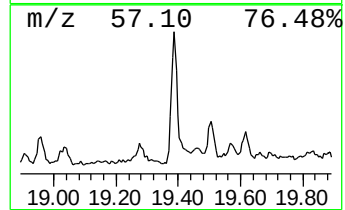
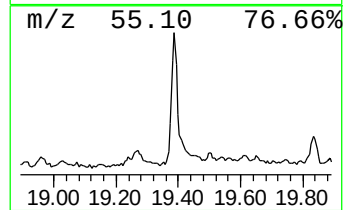
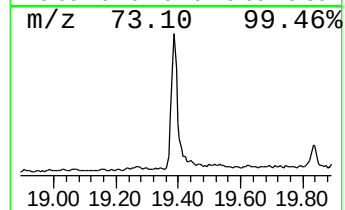
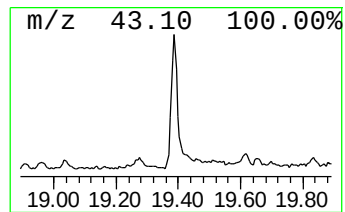
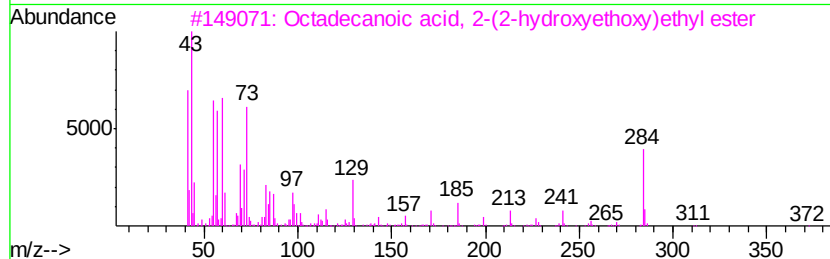
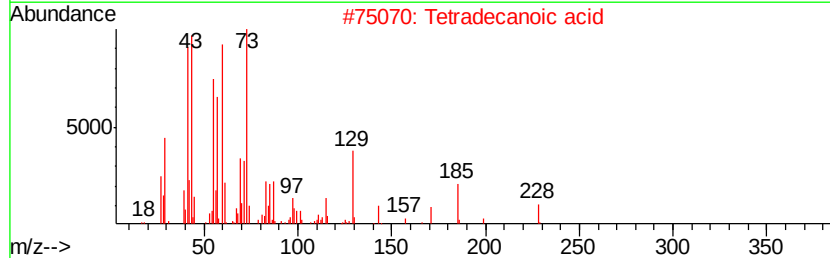
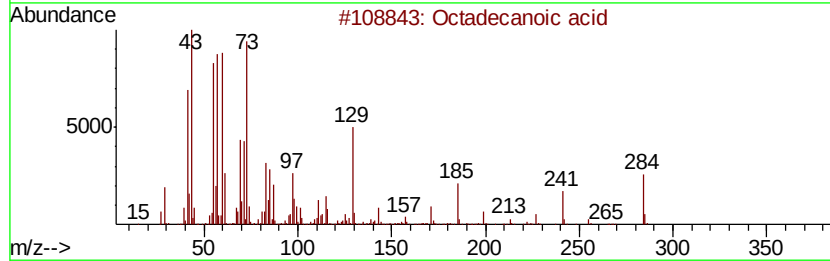
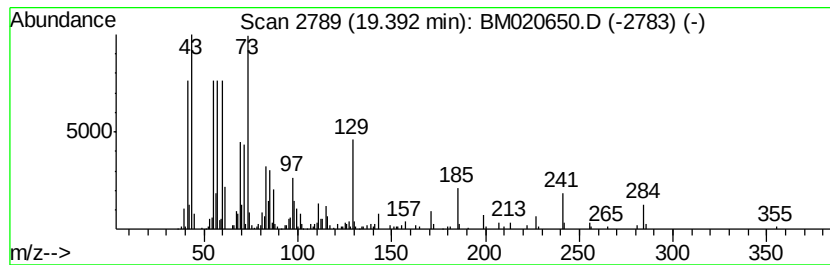
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
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Peak Number 6 Octadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.39	3.49 ng	631806	Chrysene-d12	21.39

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octadecanoic acid	284	C18H36O2	000057-11-4	99
2		Tetradecanoic acid	228	C14H28O2	000544-63-8	93
3		Octadecanoic acid, 2-(2-hydroxyve...	372	C22H44O4	000106-11-6	91
4		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	83
5		Pentadecanoic acid	242	C15H30O2	001002-84-2	72



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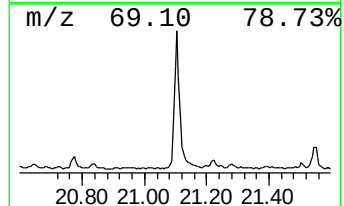
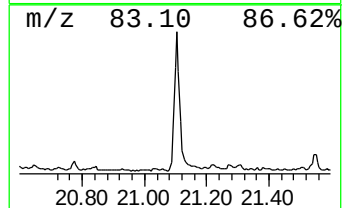
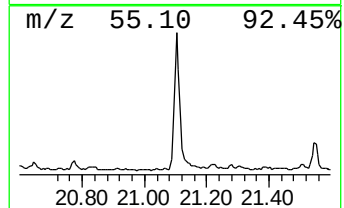
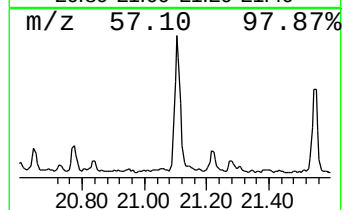
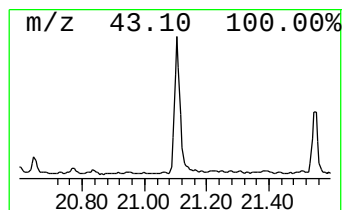
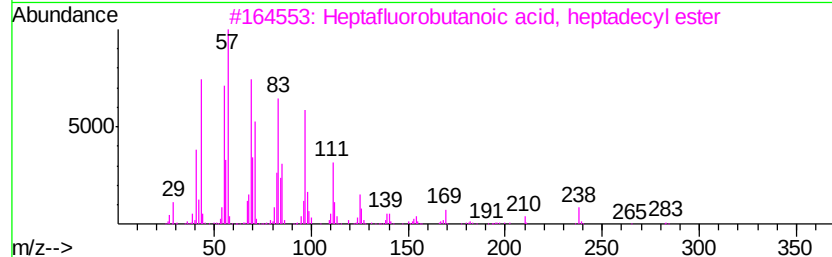
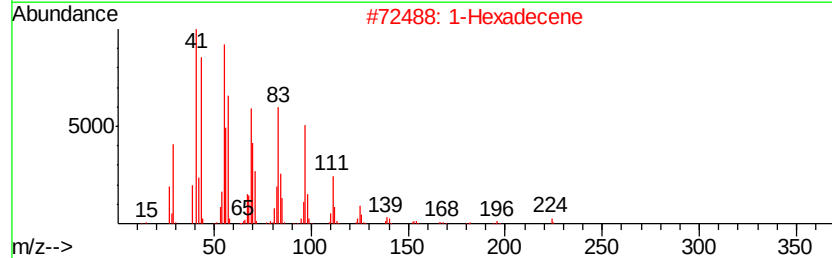
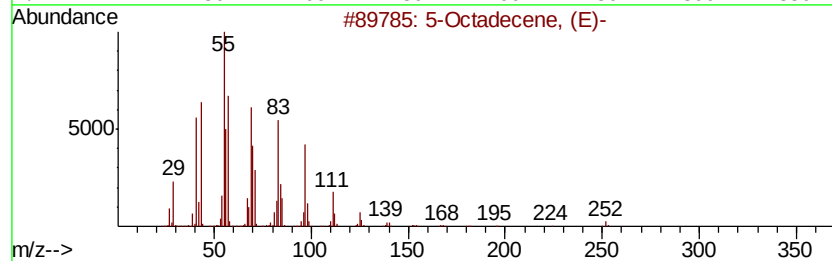
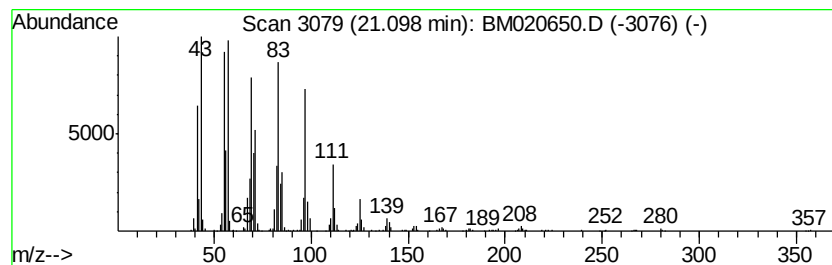
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ClientSampleId :
TEST-PIT-5

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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 5-Octadecene, (E)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.10	6.22 ng	1124110	Chrysene-d12	21.39	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Octadecene, (E)-	252	C18H36	007206-21-5	94
2	1-Hexadecene	224	C16H32	000629-73-2	94
3	Heptafluorobutanoic acid, heptad...	452	C21H35F7O2	1000282-97-3	94
4	Cyclotetracosane	336	C24H48	000297-03-0	91
5	Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM060119\
Data File : BM020650.D
Acq On : 01 Jun 2019 14:56
Operator : HP/JU
Sample : K3075-09
Misc :
ALS Vial : 8 Sample Multiplier: 1

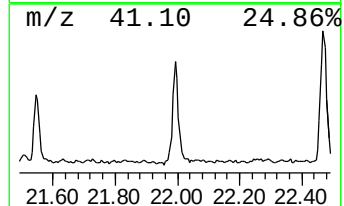
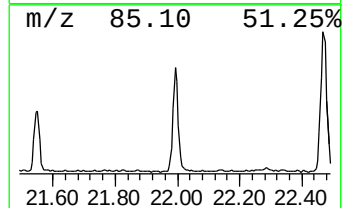
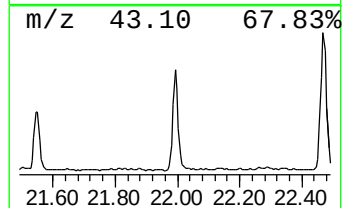
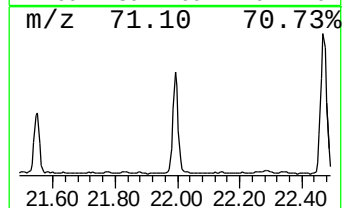
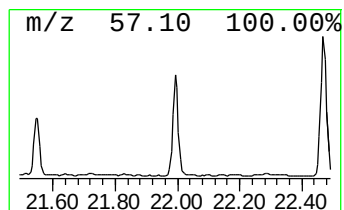
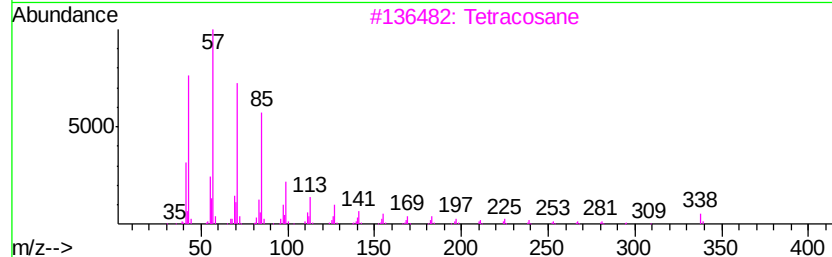
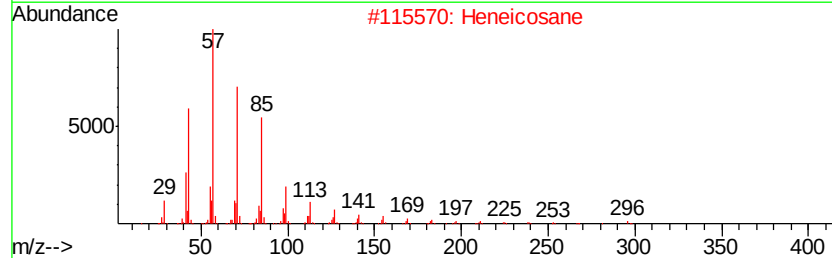
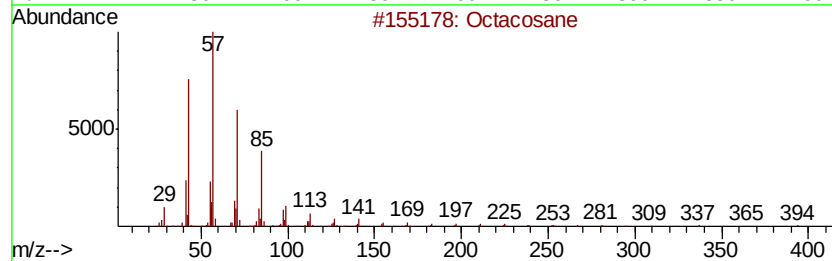
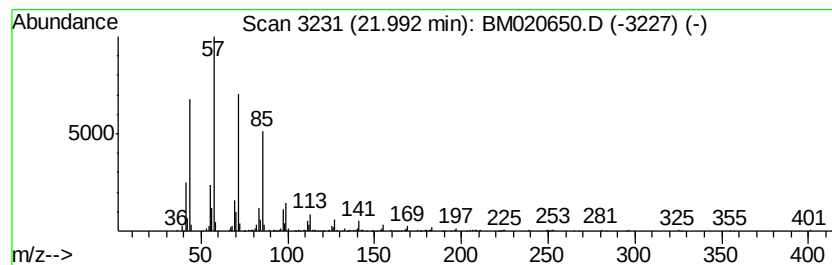
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Octacosane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.99	3.34 ng	604512	Chrysene-d12	21.39
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Octacosane	394 C28H58	000630-02-4	91
2	Heneicosane	296 C21H44	000629-94-7	91
3	Tetracosane	338 C24H50	000646-31-1	91
4	Docosane, 11-butyl-	366 C26H54	013475-76-8	91
5	Eicosane	282 C20H42	000112-95-8	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM060119\
Data File : BM020650.D
Acq On : 01 Jun 2019 14:56
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ClientSampleId :
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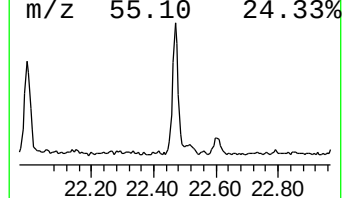
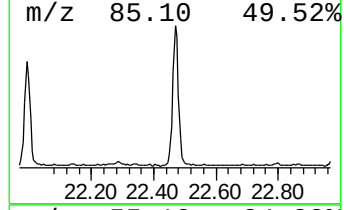
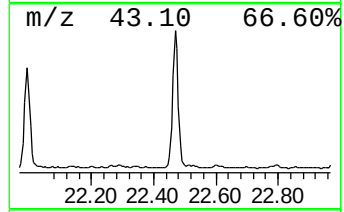
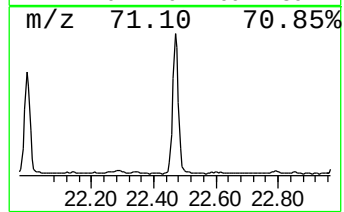
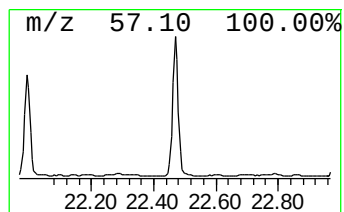
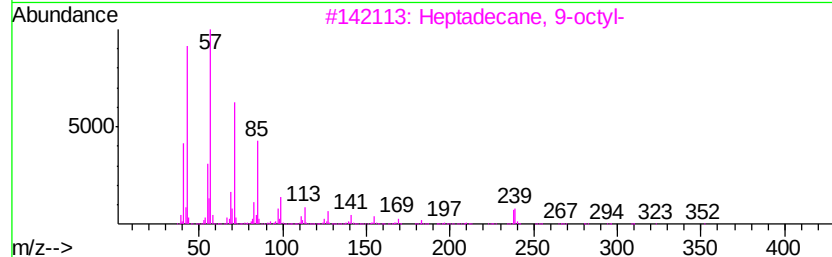
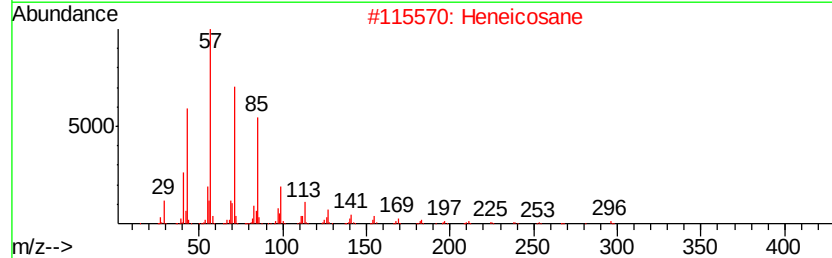
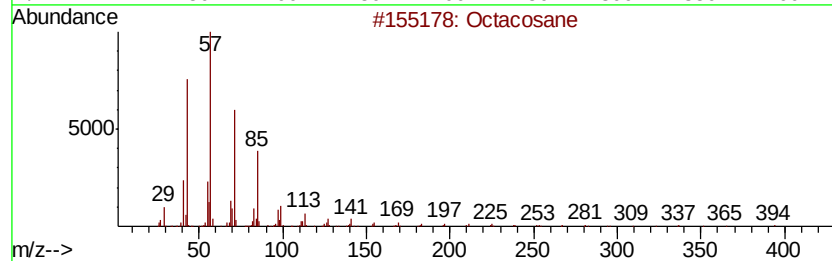
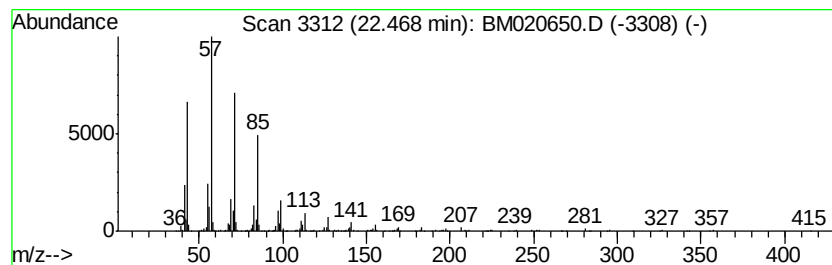
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Heneicosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.47	4.52 ng	817363	Chrysene-d12	21.39

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octacosane	394	C28H58	000630-02-4	91
2		Heneicosane	296	C21H44	000629-94-7	91
3		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
4		Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	91
5		Tetracosane	338	C24H50	000646-31-1	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM060119\
Data File : BM020650.D
Acq On : 01 Jun 2019 14:56
Operator : HP/JU
Sample : K3075-09
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
TEST-PIT-5

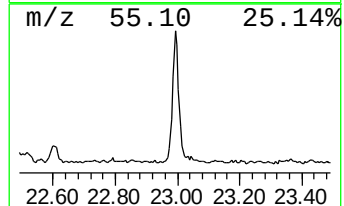
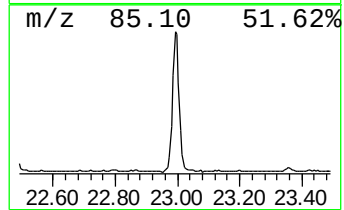
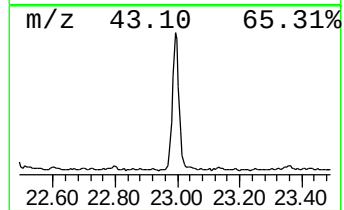
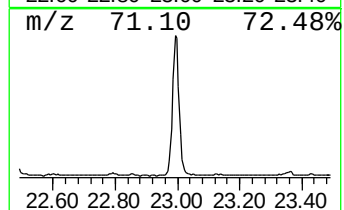
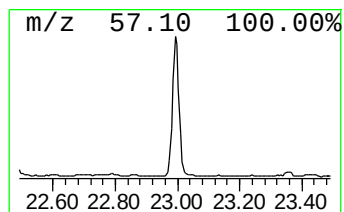
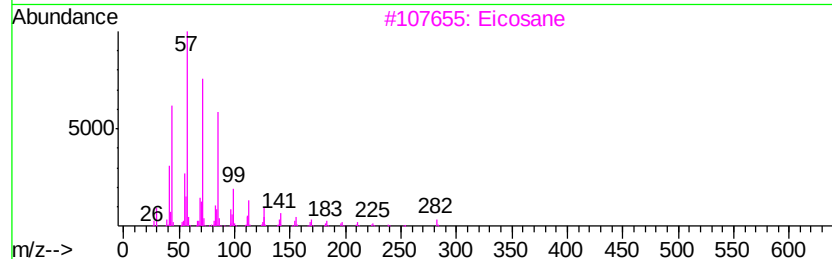
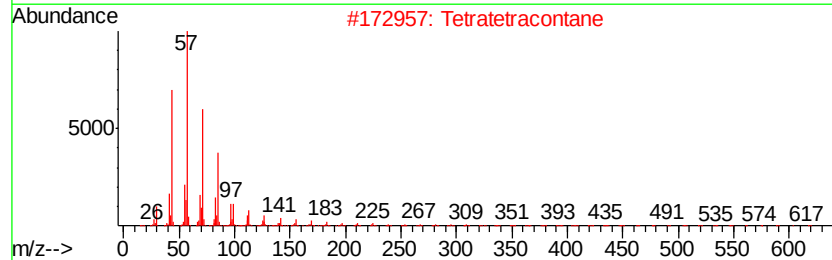
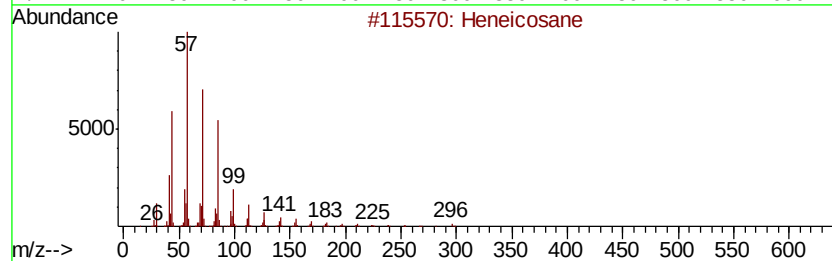
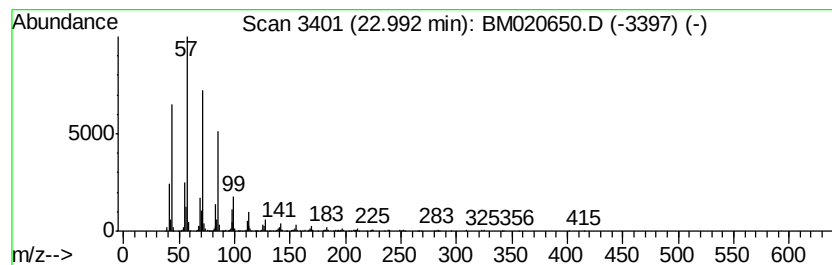
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Tetratetracontane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.99	5.18 ng	899945	Perylene-d12	23.68

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	91
2		Tetratetracontane	619	C44H90	007098-22-8	91
3		Eicosane	282	C20H42	000112-95-8	91
4		Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	91
5		Tetracosane	338	C24H50	000646-31-1	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM060119\
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Sample : K3075-09
Misc :
ALS Vial : 8 Sample Multiplier: 1

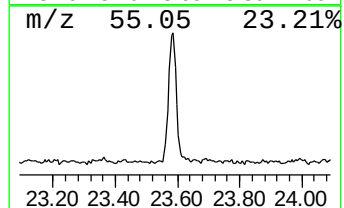
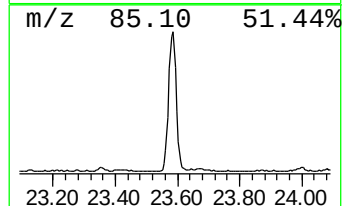
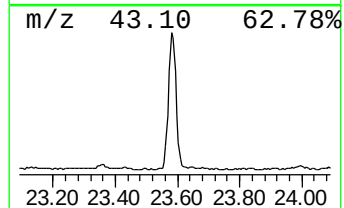
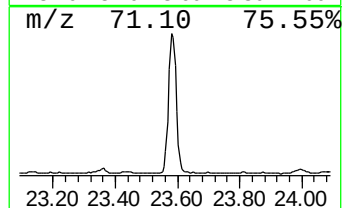
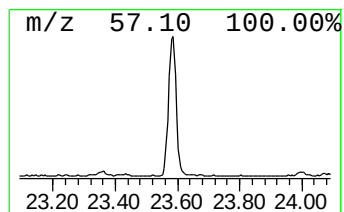
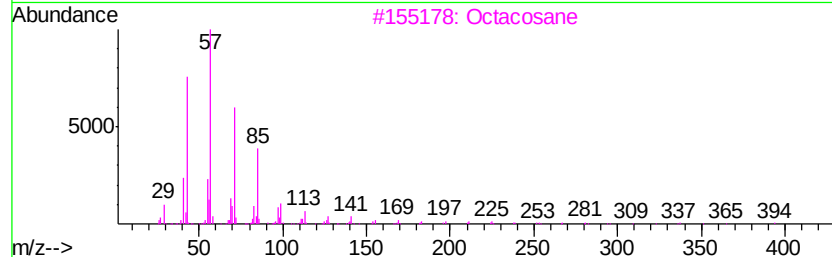
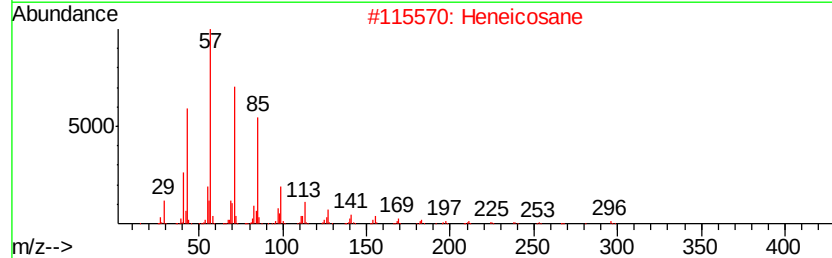
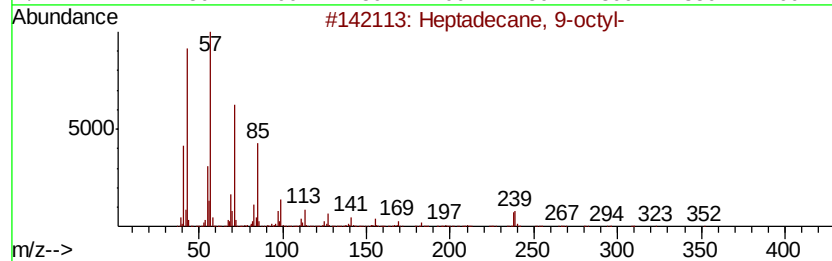
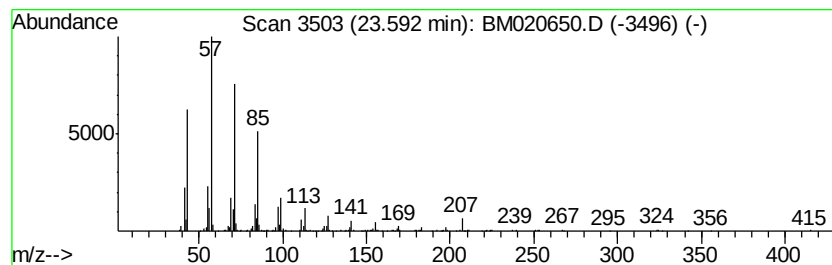
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Heptadecane, 9-octyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD		R.T.
23.59	5.58 ng	969447	Perylene-d12		23.68
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
2	Heneicosane	296	C21H44	000629-94-7	91
3	Octacosane	394	C28H58	000630-02-4	91
4	Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	91
5	triacontane	422	C30H62	000638-68-6	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM060119\
Data File : BM020650.D
Acq On : 01 Jun 2019 14:56
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Misc :
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ClientSampleId :
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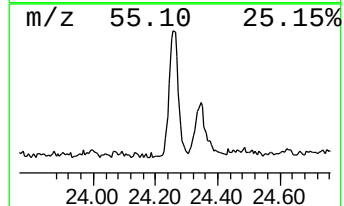
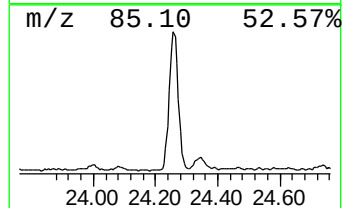
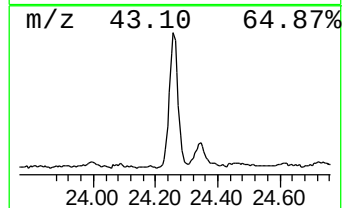
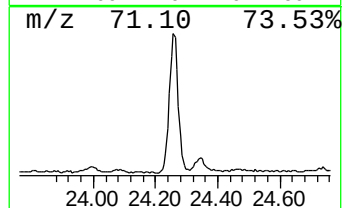
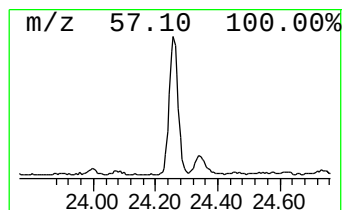
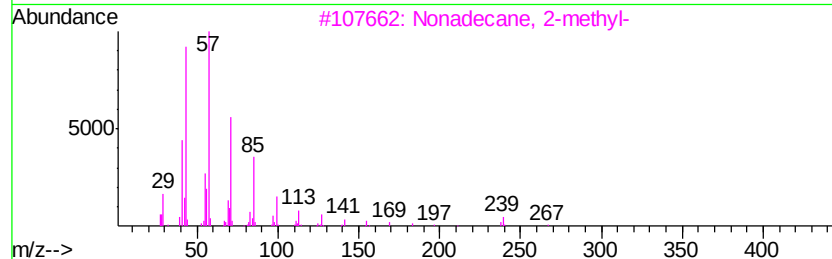
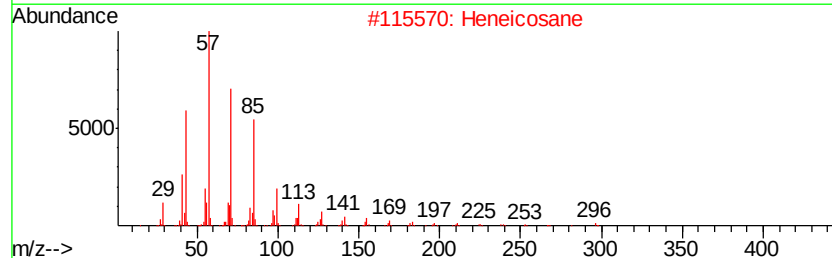
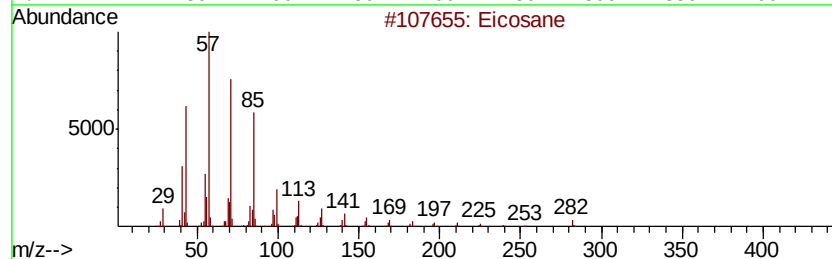
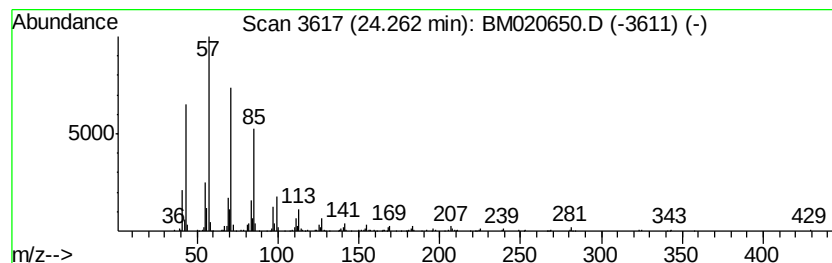
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Eicosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.26	4.27 ng	741617	Perylene-d12	23.68

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	94
2		Heneicosane	296	C21H44	000629-94-7	90
3		Nonadecane, 2-methyl-	282	C20H42	001560-86-7	87
4		Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	87
5		Eicosane, 2-methyl-	296	C21H44	001560-84-5	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM060119\
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Acq On : 01 Jun 2019 14:56
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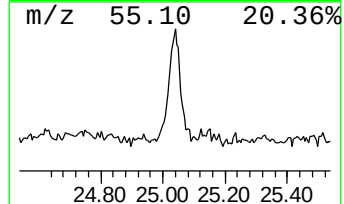
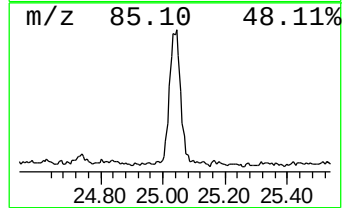
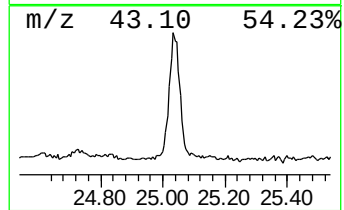
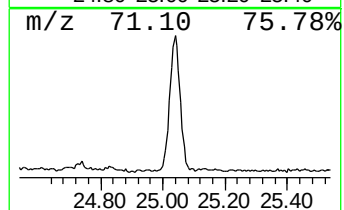
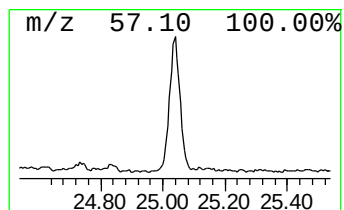
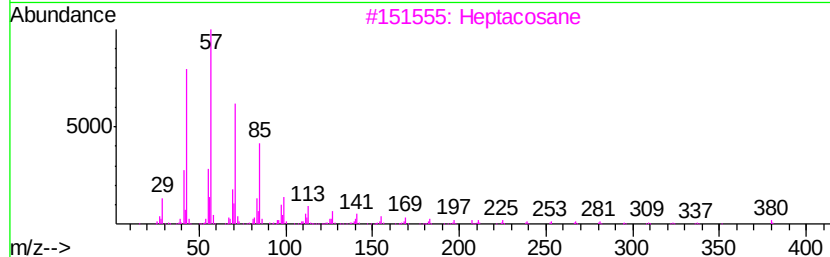
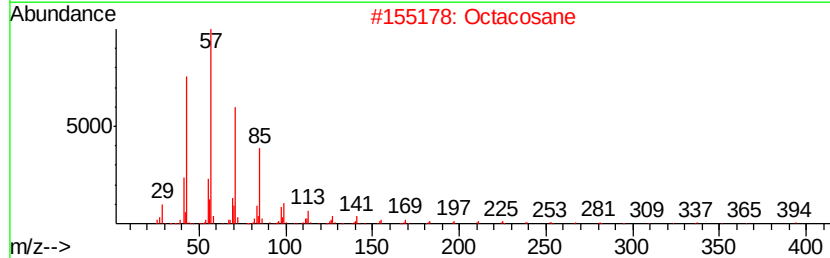
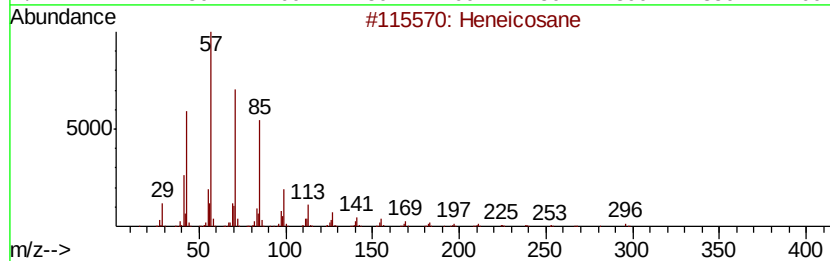
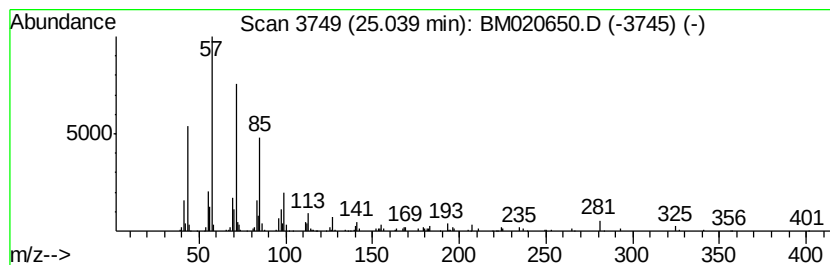
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.04	2.40 ng	416638	Perylene-d12	23.68

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	90
2		Octacosane	394	C28H58	000630-02-4	87
3		Heptacosane	380	C27H56	000593-49-7	87
4		Nonadecane, 2-methyl-	282	C20H42	001560-86-7	86
5		Docosane, 11-butyl-	366	C26H54	013475-76-8	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM060119\
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Sample : K3075-09
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
TEST-PIT-5

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\8270-BM053119.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
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unknown7.36	7.36	107.2	ng	9175680	1	7.83	1711380	20.0
Resorcinol	11.64	3.4	ng	401113	2	10.62	2355660	20.0
n-Hexadecanoic acid	18.11	7.3	ng	1342990	4	17.21	3658670	20.0
Octadecanoic acid	19.39	3.5	ng	631806	5	21.39	3616300	20.0
5-Octadecene, (E)-	21.10	6.2	ng	1124110	5	21.39	3616300	20.0
Octacosane	21.99	3.3	ng	604512	5	21.39	3616300	20.0
Heneicosane	22.47	4.5	ng	817363	5	21.39	3616300	20.0
Tetratetracontane	22.99	5.2	ng	899945	6	23.68	3473920	20.0
Heptadecane, 9-oc...	23.59	5.6	ng	969447	6	23.68	3473920	20.0
Eicosane	24.26	4.3	ng	741617	6	23.68	3473920	20.0
Nonadecane, 2-met...	25.04	2.4	ng	416638	6	23.68	3473920	20.0