

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032119\
 Data File : BM019301.D
 Acq On : 21 Mar 2019 21:33
 Operator : JU/SJ
 Sample : K2004-04
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SB-19(13-14)

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-BM031119.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	5.234	375	382	399	rBV	4328942	6194580	62.87%	7.382%
2	6.816	644	651	668	rBV	4722543	7585144	76.98%	9.039%
3	7.163	703	710	728	rBV	4811865	7498620	76.11%	8.936%
4	7.628	783	789	796	rBV	867338	1333108	13.53%	1.589%
5	7.946	835	843	854	rBV	3250838	5093391	51.69%	6.070%
6	8.798	980	988	1007	rBV	2424093	4017078	40.77%	4.787%
7	9.875	1163	1171	1190	rBV	456545	908016	9.22%	1.082%
8	10.410	1255	1262	1275	rBV	1144067	1887085	19.15%	2.249%
9	11.251	1395	1405	1416	rBV	102230	204324	2.07%	0.243%
10	12.898	1677	1685	1698	rBV	6883989	9744734	98.90%	11.613%
11	13.233	1736	1742	1747	rBV2	75553	146060	1.48%	0.174%
12	13.498	1783	1787	1794	rVB3	75821	118125	1.20%	0.141%
13	13.751	1824	1830	1844	rVB	470444	726648	7.38%	0.866%
14	14.286	1914	1921	1930	rVB2	2071072	3004315	30.49%	3.580%
15	15.786	2165	2176	2190	rBV	6226098	8979426	91.14%	10.701%
16	15.892	2191	2194	2201	rVB2	61327	114754	1.16%	0.137%
17	17.045	2383	2390	2398	rBV2	2208091	3206730	32.55%	3.821%
18	17.939	2536	2542	2551	rBV	989720	1374616	13.95%	1.638%
19	18.421	2619	2624	2632	rBV	233497	330581	3.36%	0.394%
20	18.610	2651	2656	2663	rBV	148306	321243	3.26%	0.383%
21	19.221	2756	2760	2768	rBV	280021	479271	4.86%	0.571%
22	19.686	2833	2839	2844	rBV	8201972	9852835	100.00%	11.742%
23	19.980	2887	2889	2893	rVB	184931	180544	1.83%	0.215%
24	20.951	3050	3054	3064	rVB	1108735	1763683	17.90%	2.102%
25	21.245	3100	3104	3109	rBV	2088114	2539966	25.78%	3.027%
26	21.756	3188	3191	3195	rBV5	265164	314961	3.20%	0.375%
27	21.809	3198	3200	3203	rVV	230766	272036	2.76%	0.324%
28	21.845	3203	3206	3214	rVB2	412518	720924	7.32%	0.859%
29	22.768	3360	3363	3368	rVB	313589	376198	3.82%	0.448%
30	23.327	3455	3458	3463	rVB3	167039	249647	2.53%	0.298%
31	23.474	3477	3483	3490	rVB2	1442694	2711099	27.52%	3.231%
32	24.074	3581	3585	3593	rBV5	198064	439894	4.46%	0.524%
33	27.368	4141	4145	4156	rVB6	412125	1224953	12.43%	1.460%

LSC Area Percent Report

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SB-19(13-14)

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

Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\8270-BM031119.M
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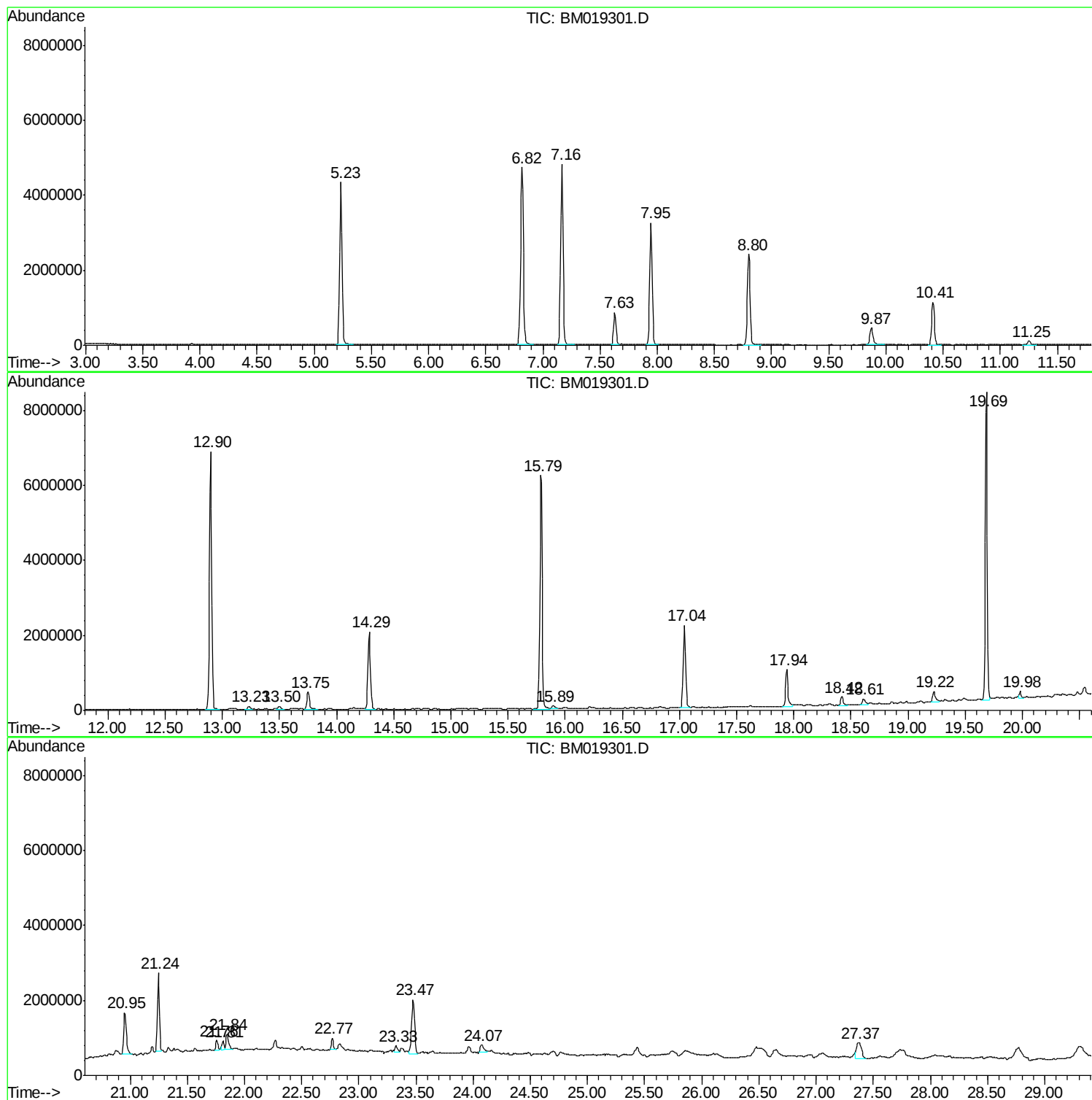
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM031119.M
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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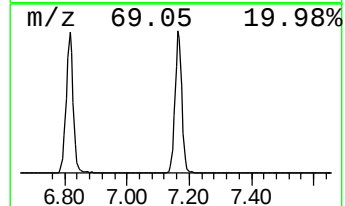
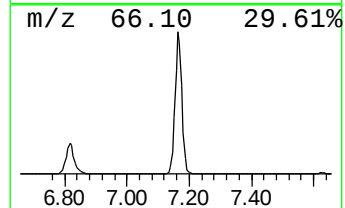
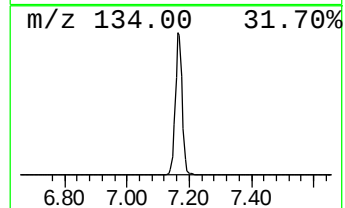
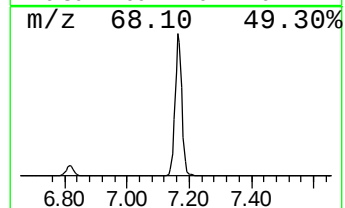
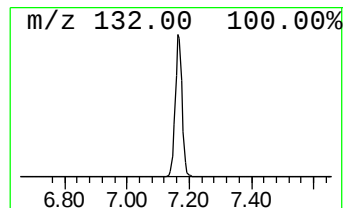
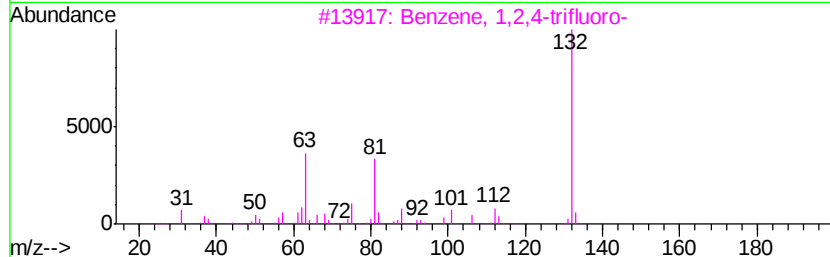
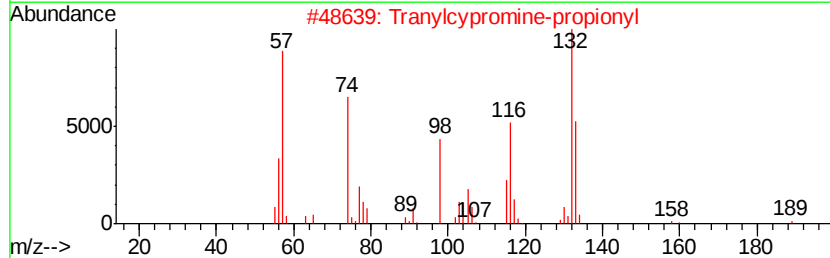
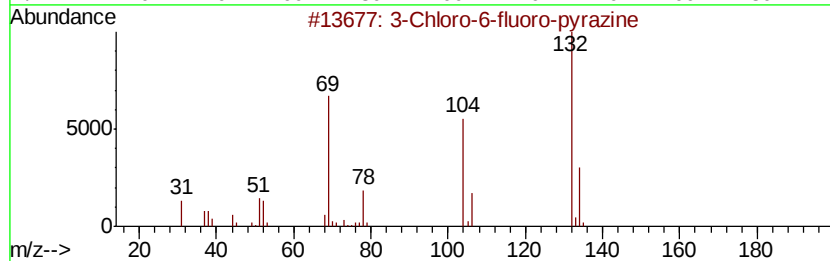
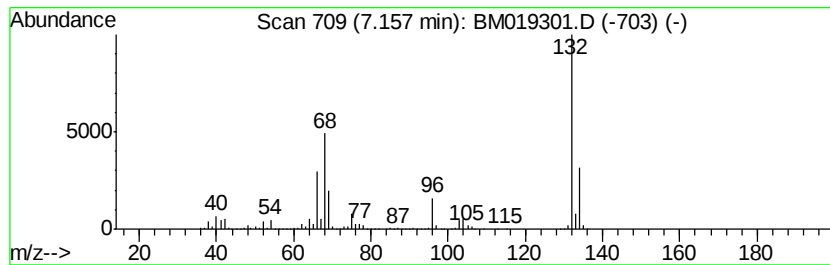
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 unknown7.16 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.16	112.50 ng	7498620	1,4-Dichlorobenzene-d4	7.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
2		Tranvlcypropamine-propionyl	189	C12H15NO	1000123-86-3	10
3		Benzene, 1,2,4-trifluoro-	132	C6H3F3	000367-23-7	9
4		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9
5		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9



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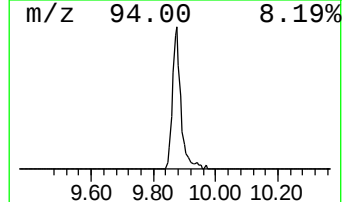
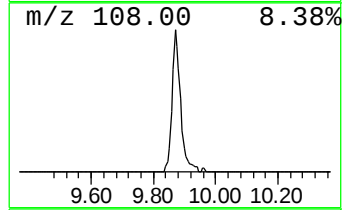
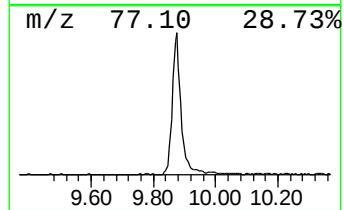
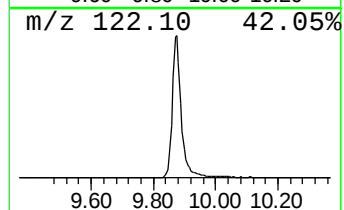
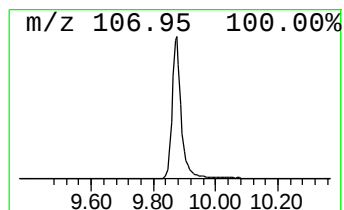
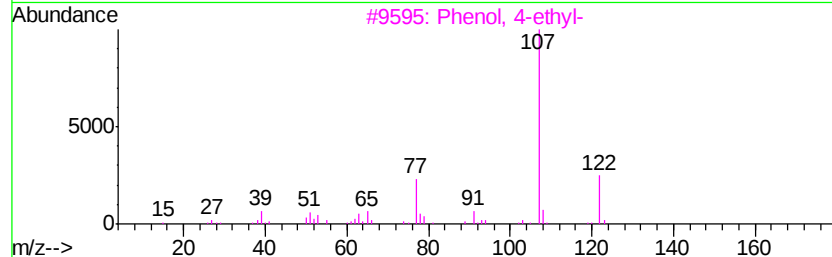
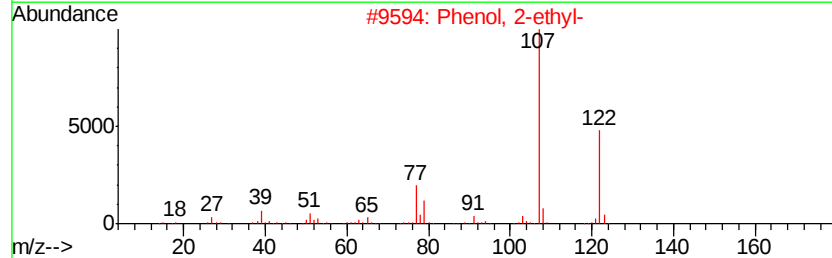
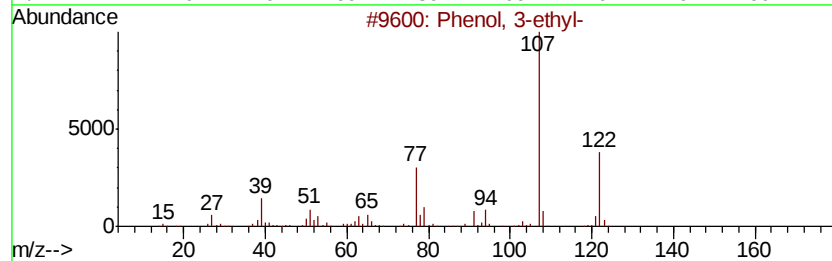
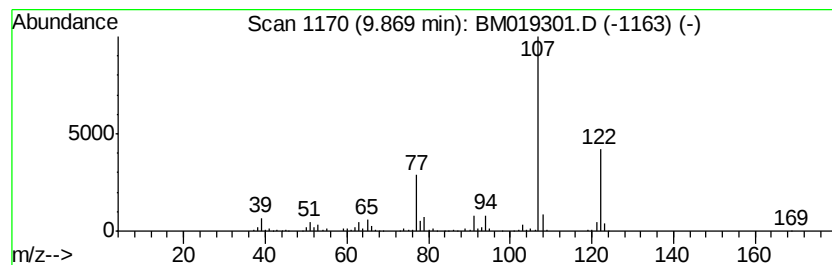
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 3 Phenol, 3-ethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.87	9.62 ng	908016	Naphthalene-d8	10.41

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenol, 3-ethyl-	122	C8H10O	000620-17-7	94
2		Phenol, 2-ethyl-	122	C8H10O	000090-00-6	91
3		Phenol, 4-ethyl-	122	C8H10O	000123-07-9	86
4		3,3-Dimethyl-6-methylenecyclohexene	122	C9H14	020185-16-4	72
5		2-Isopropylpyrazine	122	C7H10N2	029460-90-0	59



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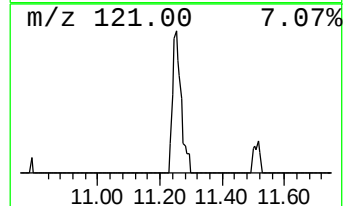
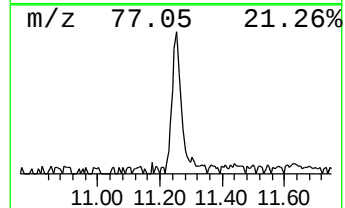
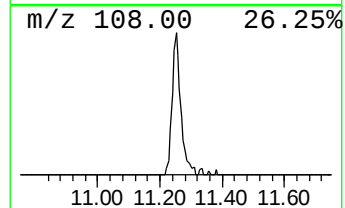
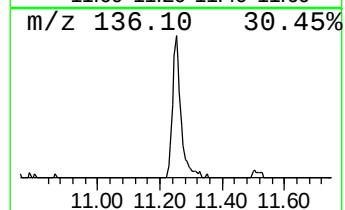
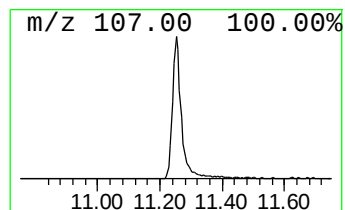
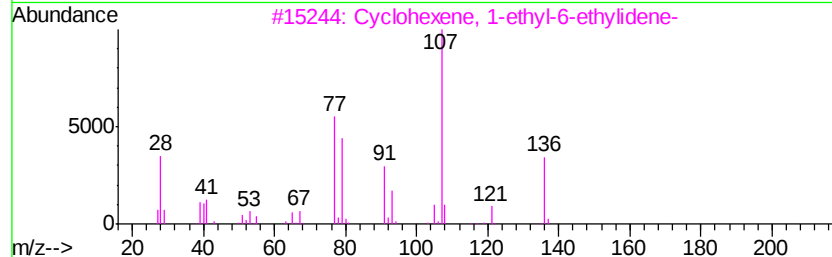
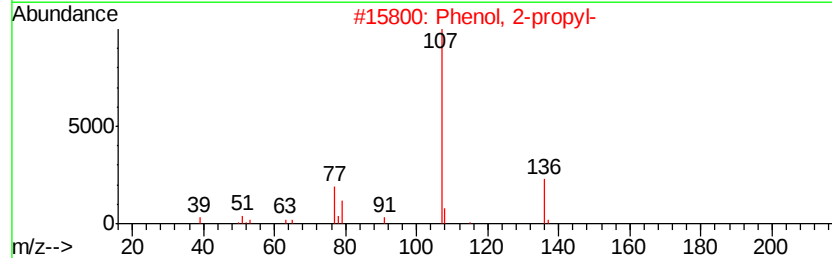
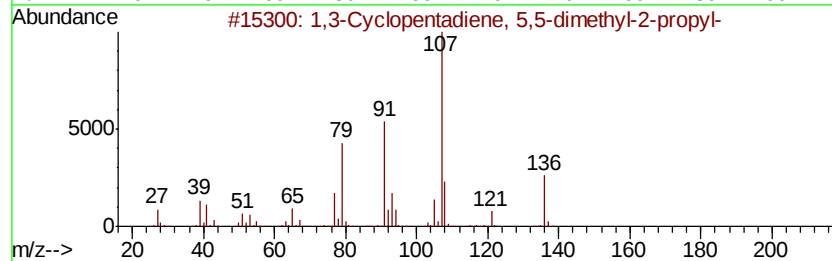
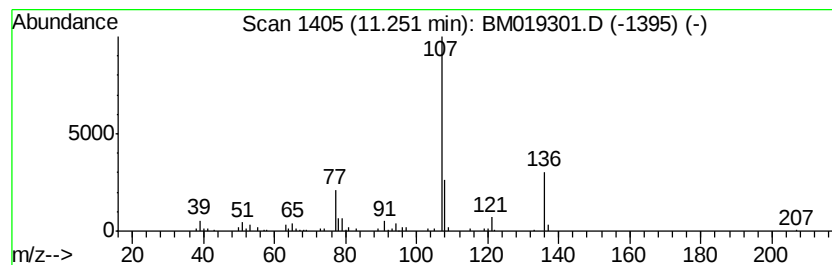
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 4 1,3-Cyclopentadiene, 5,5-di... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.25	2.17 ng	204324	Naphthalene-d8	10.41

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,3-Cyclopentadiene, 5,5-dimethy...	136	C10H16	1000163-57-0	86
2		Phenol, 2-propyl-	136	C9H12O	000644-35-9	81
3		Cyclohexene, 1-ethyl-6-ethylidene-	136	C10H16	061141-57-9	59
4		Phenol, 4-propyl-	136	C9H12O	000645-56-7	58
5		Phenol, 2-butyl-	150	C10H14O	003180-09-4	53



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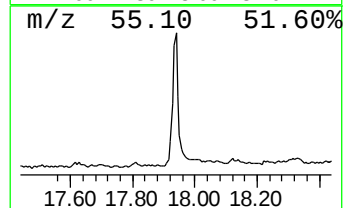
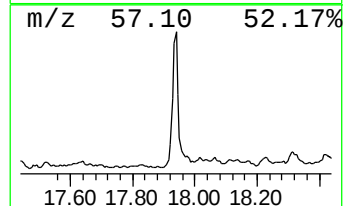
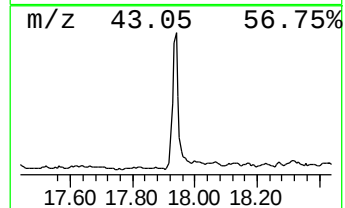
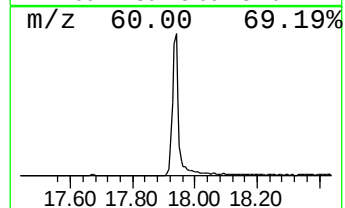
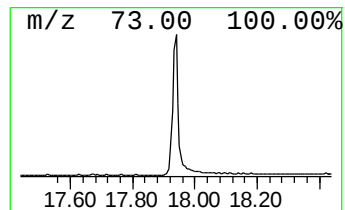
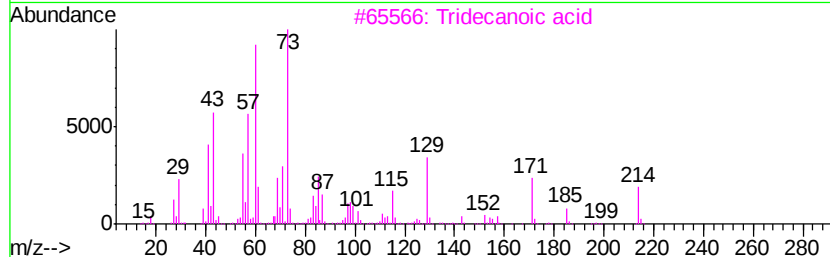
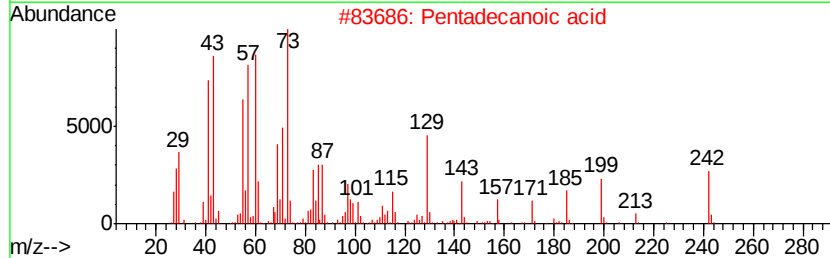
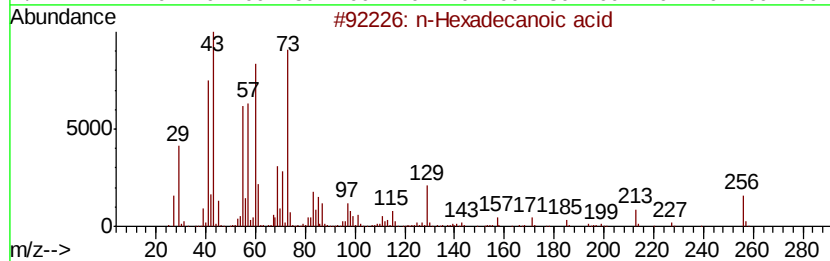
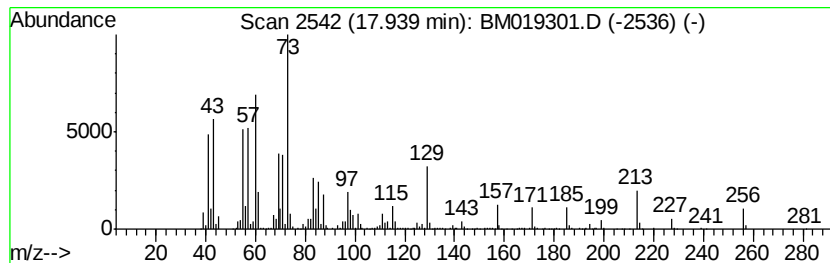
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 5 n-Hexadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.94	8.57 ng	1374620	Phenanthrene-d10	17.04

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
2	Pentadecanoic acid	242	C15H30O2	001002-84-2	62
3	Tridecanoic acid	214	C13H26O2	000638-53-9	53
4	Tetradecanoic acid	228	C14H28O2	000544-63-8	53
5	11-Bromoundecanoic acid	264	C11H21BrO2	002834-05-1	38



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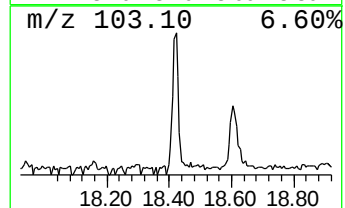
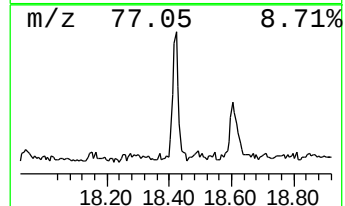
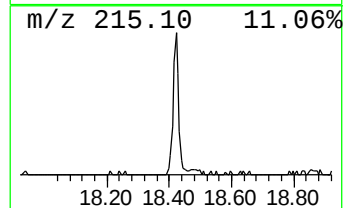
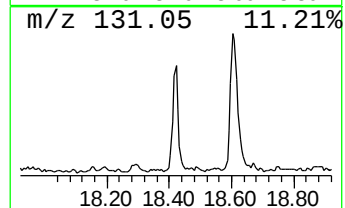
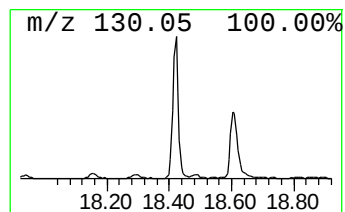
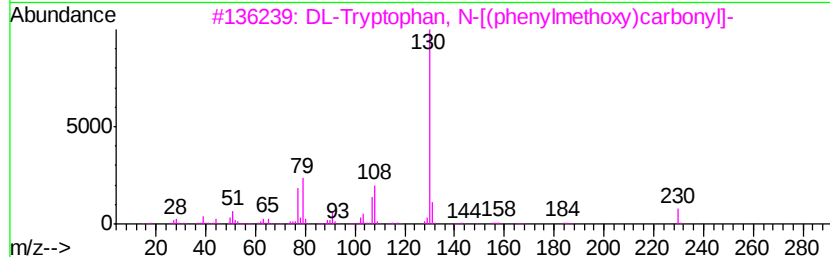
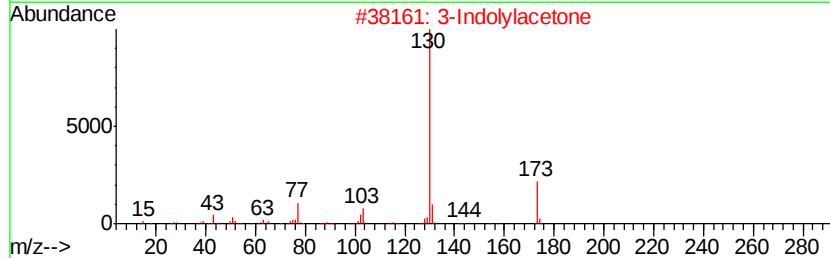
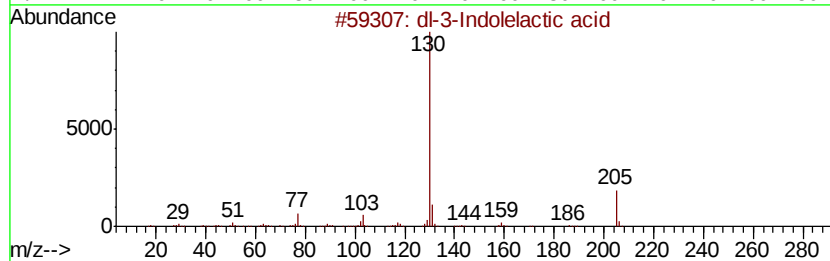
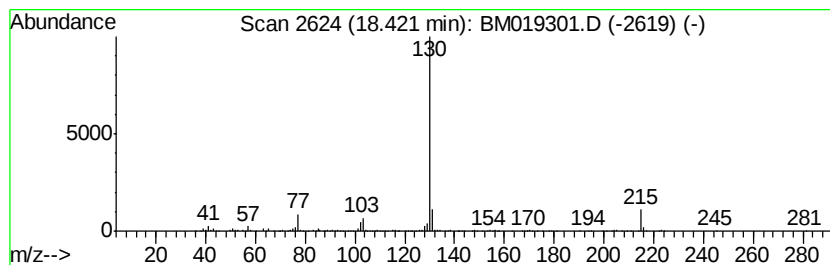
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Peak Number 6 dl-3-Indolelactic acid Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.42	2.06 ng	330581	Phenanthrene-d10	17.04

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		dl-3-Indolelactic acid	205	C11H11N03	000832-97-3	64
2		3-Indolylacetone	173	C11H11N0	001201-26-9	64
3		DL-Tryptophan, N-[(phenylmethoxyv...	338	C19H18N2O4	013058-16-7	64
4		D-Tryptophan	204	C11H12N2O2	000153-94-6	64
5		DL-Tryptophan	204	C11H12N2O2	000054-12-6	64



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Operator : JU/SJ
Sample : K2004-04
Misc :
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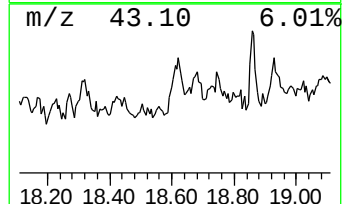
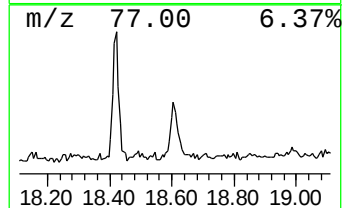
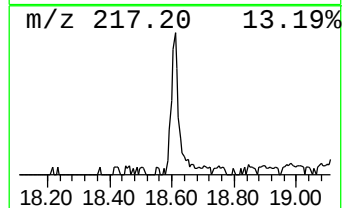
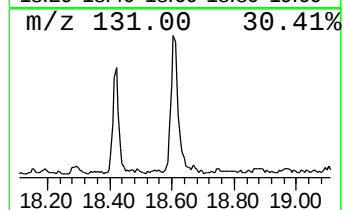
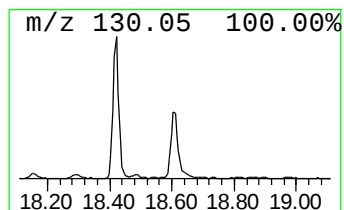
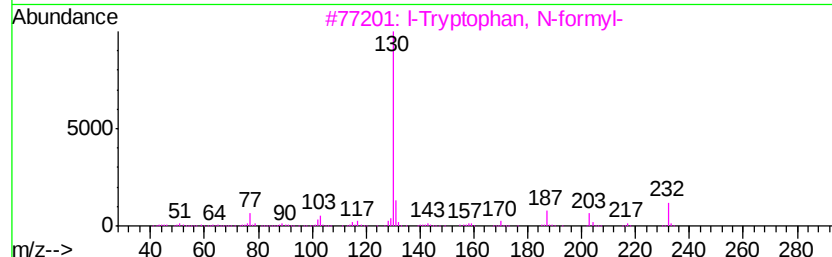
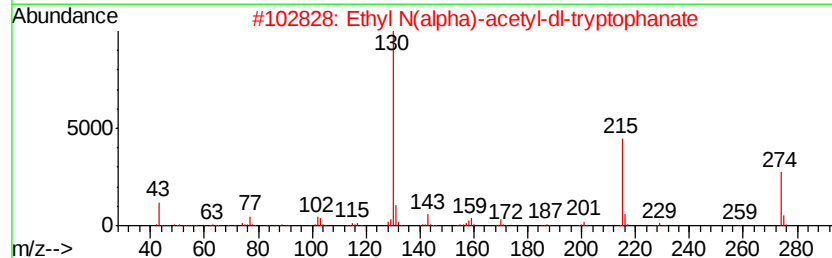
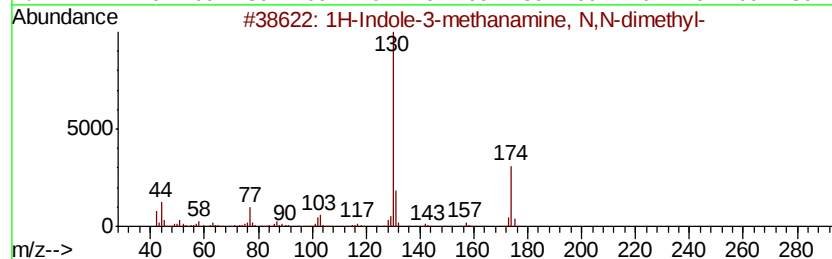
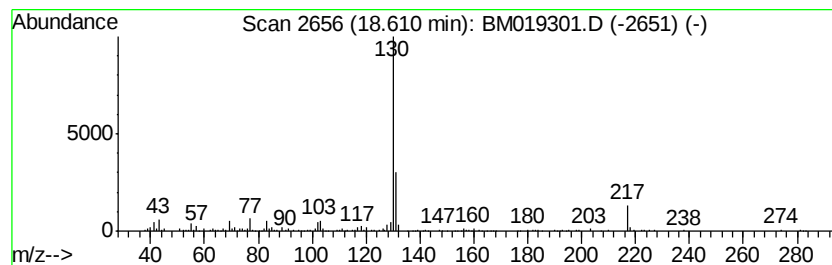
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ClientSampleId :
SB-19(13-14)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM031119.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-Indole-3-methanamine, N,... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.61	2.00 ng	321243	Phenanthrene-d10	17.04	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indole-3-methanamine, N,N-dim...	174	C11H14N2	000087-52-5	64
2	Ethyl N(alpha)-acetyl-dl-tryptop...	274	C15H18N2O3	042717-06-6	64
3	l-Tryptophan, N-formyl-	232	C12H12N2O3	1000131-24-5	64
4	Glycyl-L-tryptophan	261	C13H15N3O3	002390-74-1	59
5	L-Tryptophan, ethyl ester	232	C13H16N2O2	007479-05-2	59



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032119\
Data File : BM019301.D
Acq On : 21 Mar 2019 21:33
Operator : JU/SJ
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Misc :
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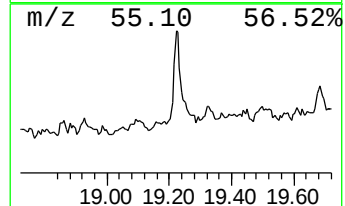
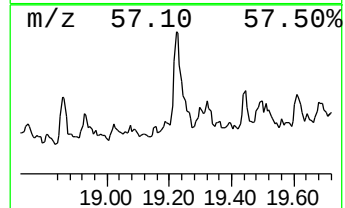
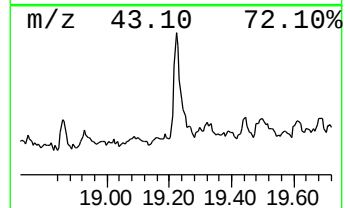
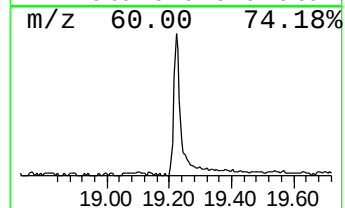
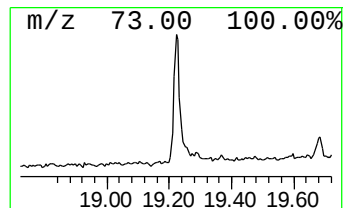
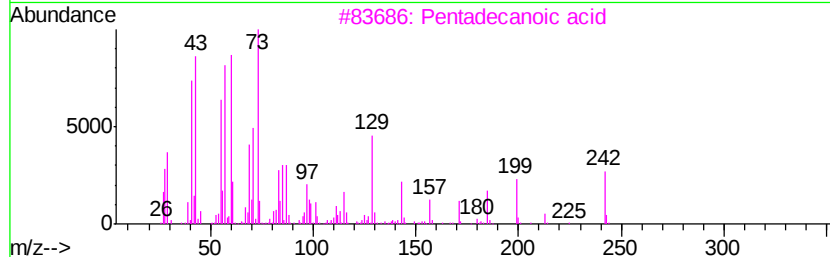
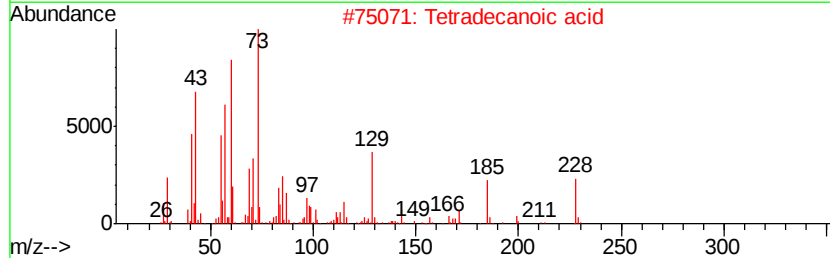
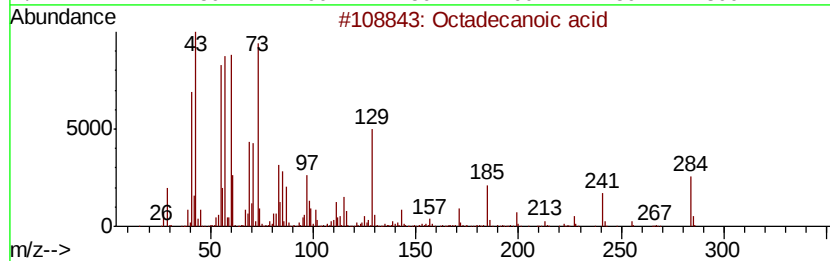
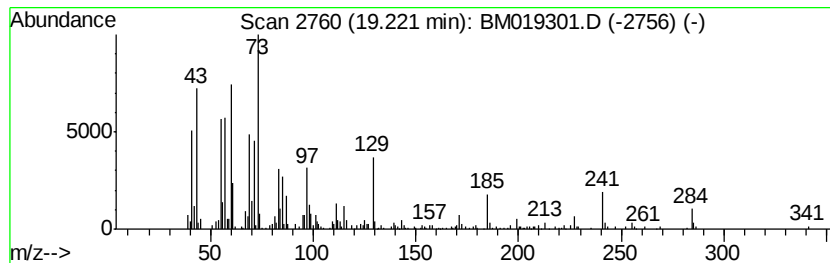
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Octadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.22	3.77 ng	479271	Chrysene-d12	21.24

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octadecanoic acid	284	C18H36O2	000057-11-4	98
2		Tetradecanoic acid	228	C14H28O2	000544-63-8	90
3		Pentadecanoic acid	242	C15H30O2	001002-84-2	81
4		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	74
5		Eicosane	282	C20H42	000112-95-8	53



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032119\
Data File : BM019301.D
Acq On : 21 Mar 2019 21:33
Operator : JU/SJ
Sample : K2004-04
Misc :
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ClientSampleId :
SB-19(13-14)

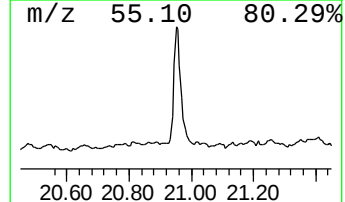
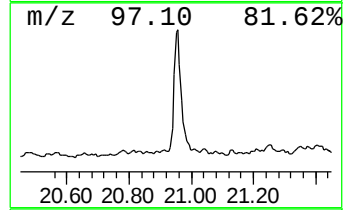
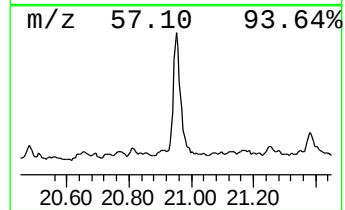
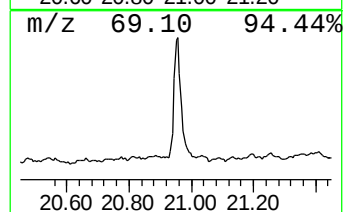
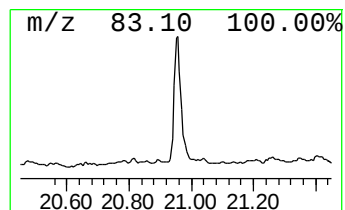
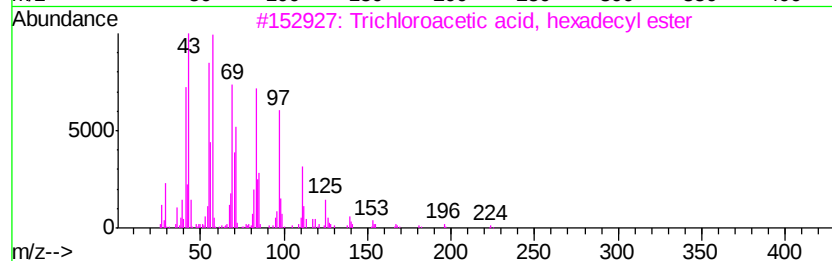
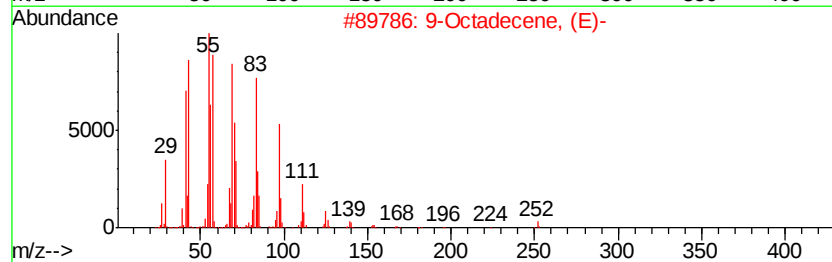
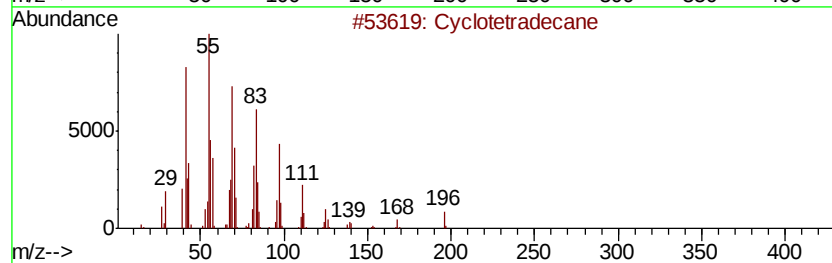
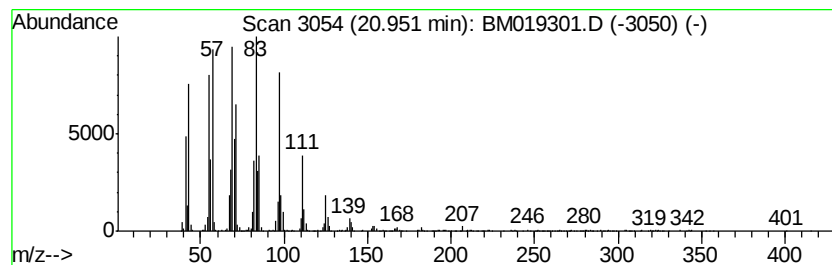
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Cyclotetradecane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.95	13.89 ng	1763680	Chrysene-d12	21.24

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclotetradecane	196	C14H28	000295-17-0	94
2		9-Octadecene, (E)-	252	C18H36	007206-25-9	93
3		Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	91
4		3-Eicosene, (E)-	280	C20H40	074685-33-9	91
5		Trichloroacetic acid, tetradecyl...	358	C16H29Cl3O2	074339-52-9	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032119\
Data File : BM019301.D
Acq On : 21 Mar 2019 21:33
Operator : JU/SJ
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Misc :
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ClientSampleId :
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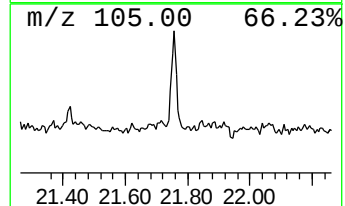
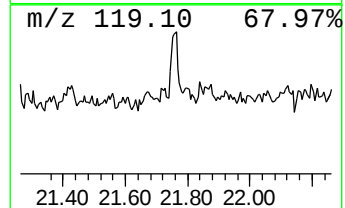
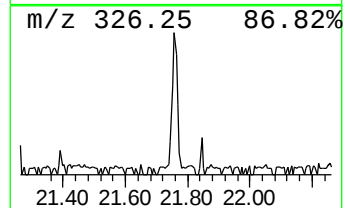
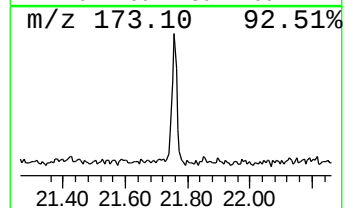
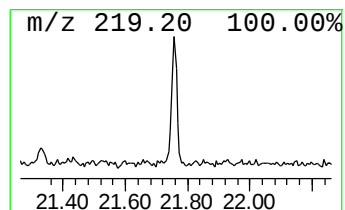
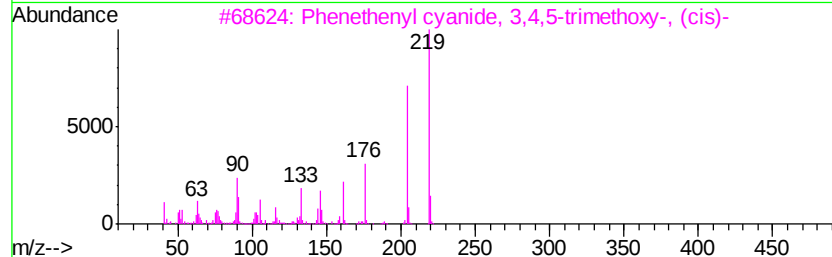
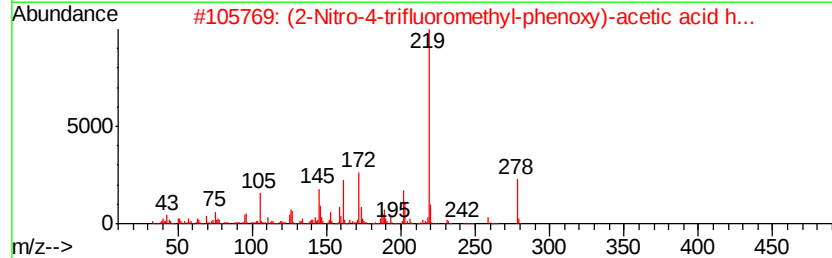
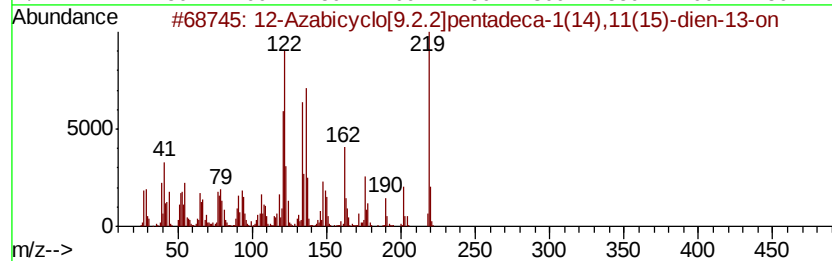
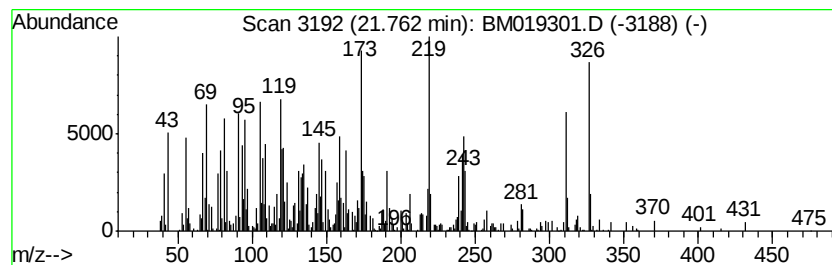
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 unknown21.76 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.76	2.48 ng	314961	Chrysene-d12	21.24

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	12-Azabicyclo[9.2.2]pentadeca-1(...	219	C14H21NO	1000185-25-1	15
2		(2-Nitro-4-trifluoromethyl-pheno...	279	C9H8F3N3O4	1000294-73-1	15
3		Phenethenyl cyanide, 3,4,5-trime...	219	C12H13N03	112369-98-9	11
4		Ethyl 3,8-dimethylimidazo[1,5-a]...	219	C11H13N3O2	1000147-83-3	11
5		Ethyl 2,3-dimethylimidazo[1,5-a]...	219	C11H13N3O2	1000147-83-4	11



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032119\
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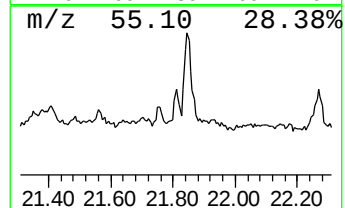
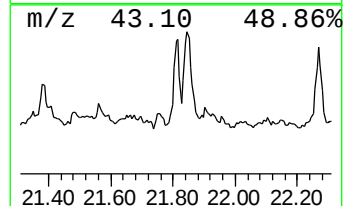
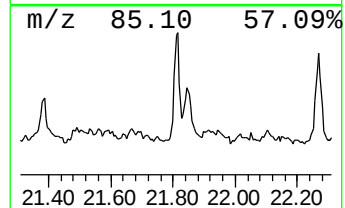
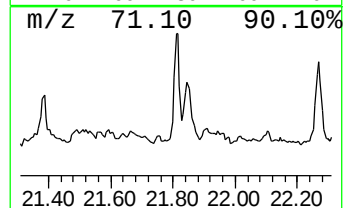
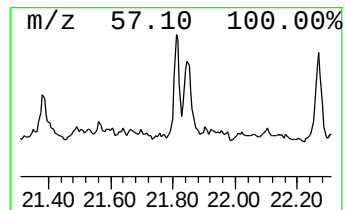
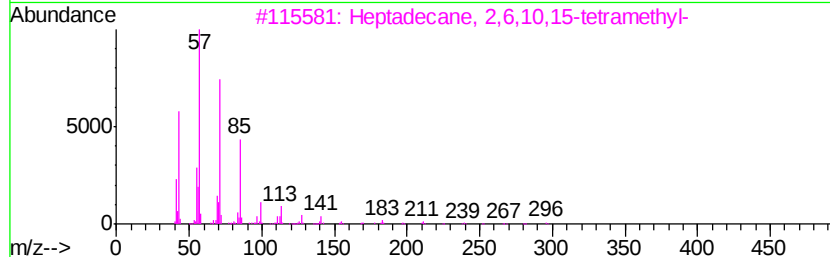
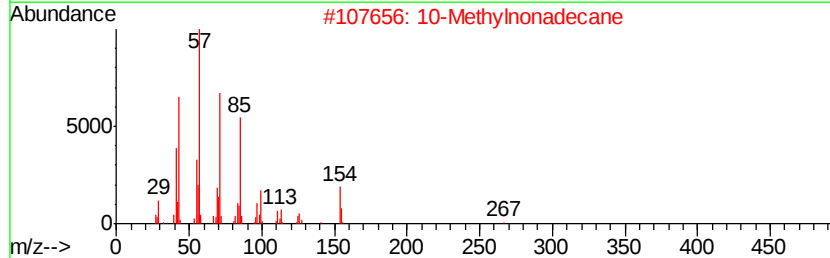
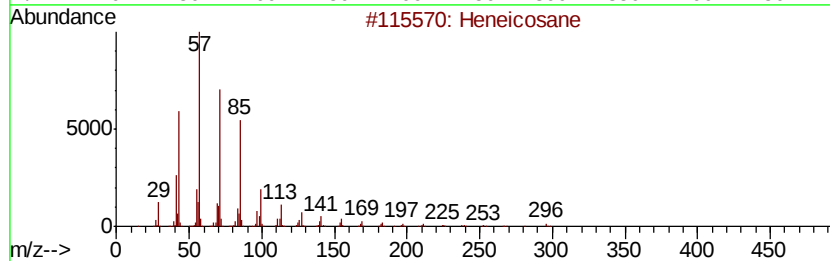
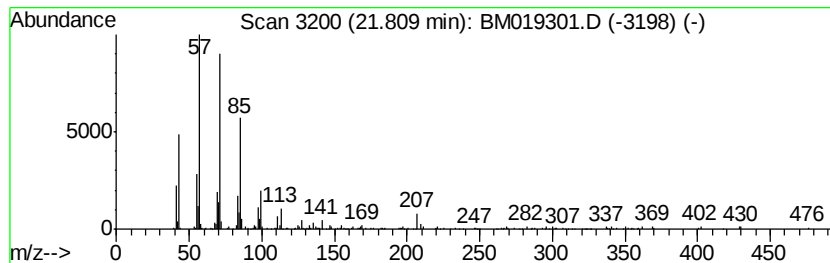
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Heneicosane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.81	2.14 ng	272036	Chrysene-d12	21.24
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Heneicosane	296 C21H44	000629-94-7	94
2	10-Methylnonadecane	282 C20H42	056862-62-5	90
3	Heptadecane, 2,6,10,15-tetramethyl-	296 C21H44	054833-48-6	87
4	Eicosane	282 C20H42	000112-95-8	78
5	2-Thiopheneacetic acid, 4-tetrad...	338 C20H34O2S	1000280-67-0	78



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032119\
Data File : BM019301.D
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ClientSampleId :
SB-19(13-14)

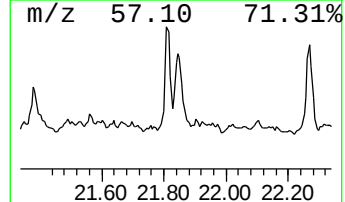
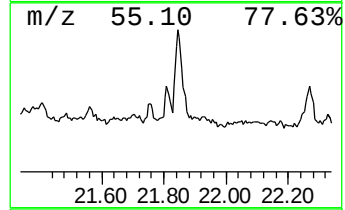
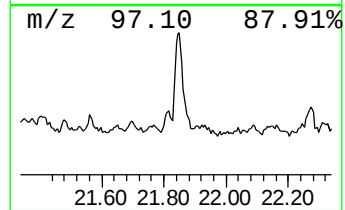
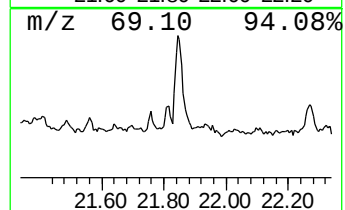
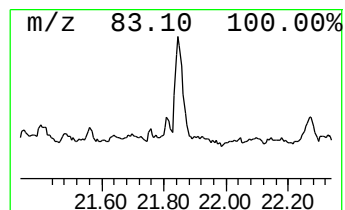
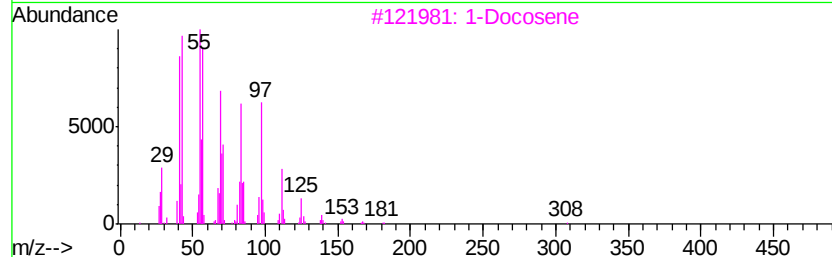
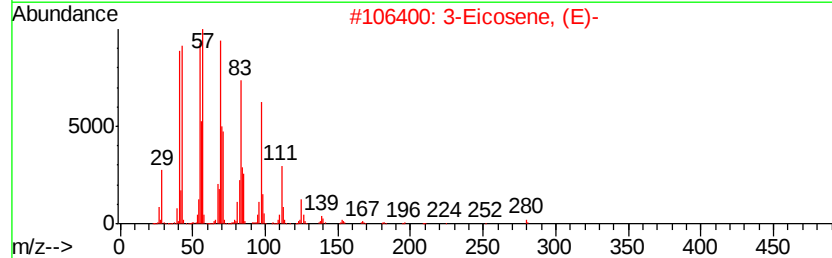
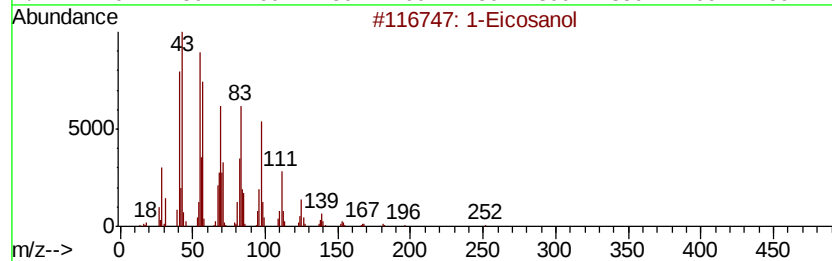
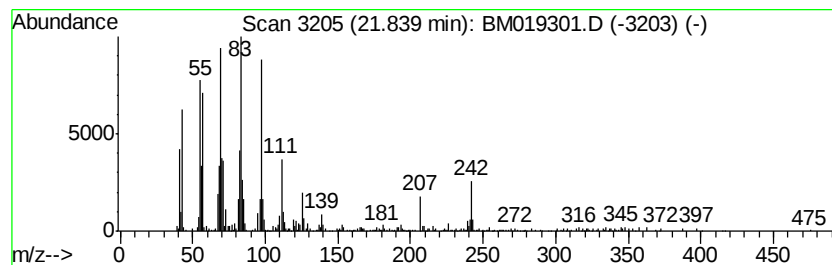
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 1-Eicosanol Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.84	5.68 ng	720924	Chrysene-d12	21.24

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Eicosanol	298	C20H42O	000629-96-9	76
2			3-Eicosene, (E)-	280	C20H40	074685-33-9	68
3			1-Docosene	308	C22H44	001599-67-3	60
4			E-15-Heptadecenal	252	C17H32O	1000130-97-9	55
5			5-Eicosene, (E)-	280	C20H40	074685-30-6	55



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032119\
Data File : BM019301.D
Acq On : 21 Mar 2019 21:33
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Sample : K2004-04
Misc :
ALS Vial : 13 Sample Multiplier: 1

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ClientSampleId :
SB-19(13-14)

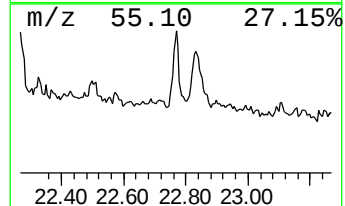
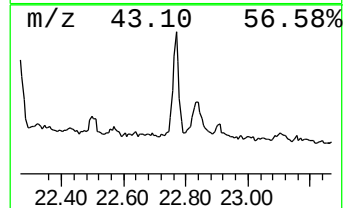
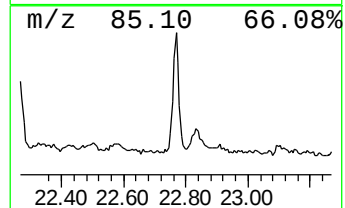
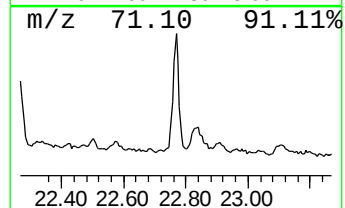
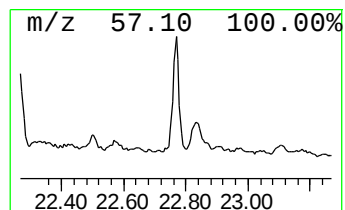
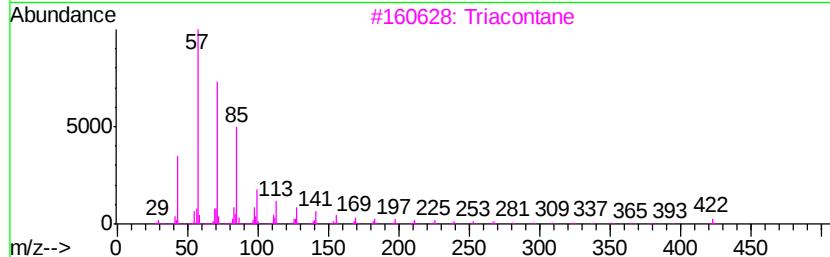
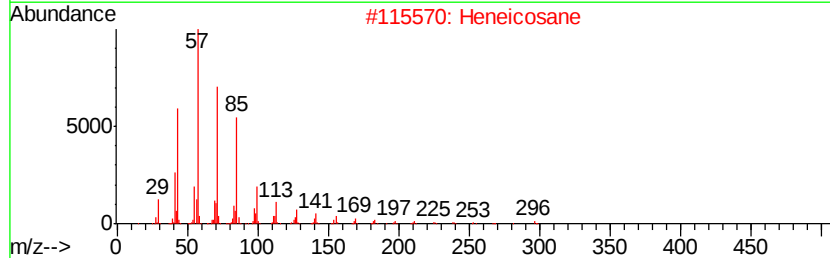
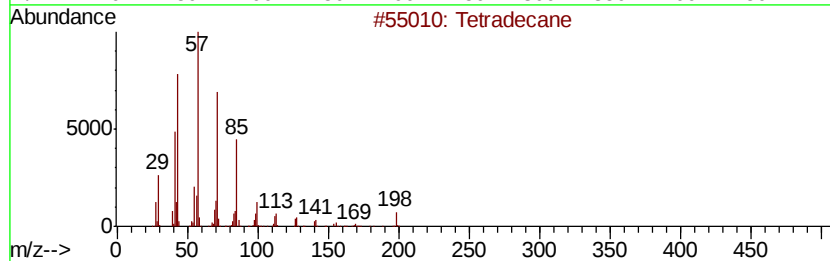
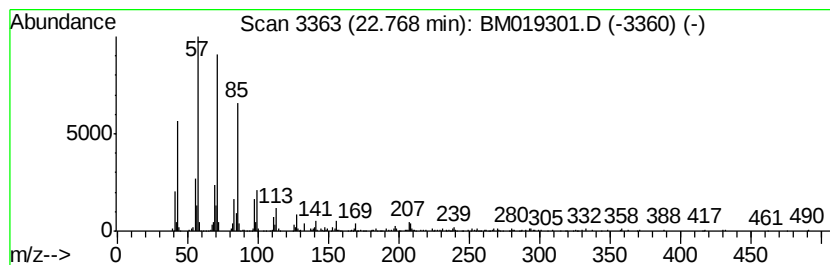
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Tetradecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.77	2.78 ng	376198	Perylene-d12	23.47

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetradecane	198	C14H30	000629-59-4	93
2		Heneicosane	296	C21H44	000629-94-7	91
3		Triacontane	422	C30H62	000638-68-6	91
4		Nonacosane	408	C29H60	000630-03-5	86
5		Docosane, 5-butyl-	366	C26H54	055282-16-1	80



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032119\
 Data File : BM019301.D
 Acq On : 21 Mar 2019 21:33
 Operator : JU/SJ
 Sample : K2004-04
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SB-19(13-14)

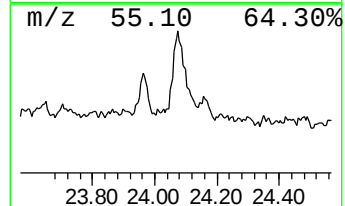
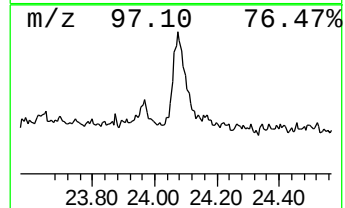
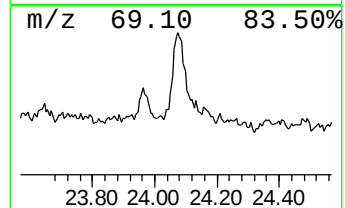
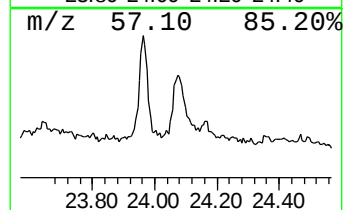
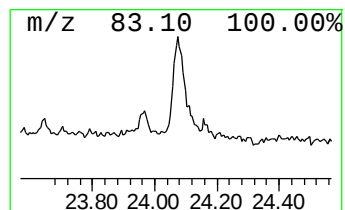
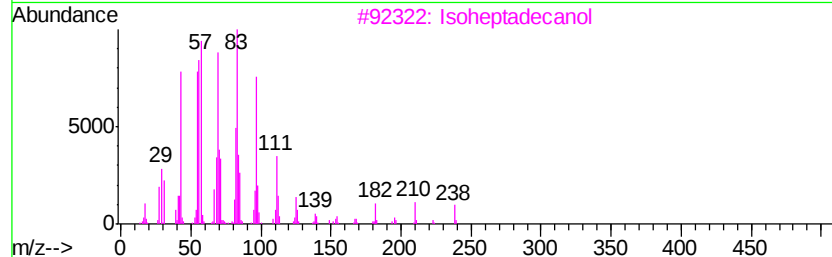
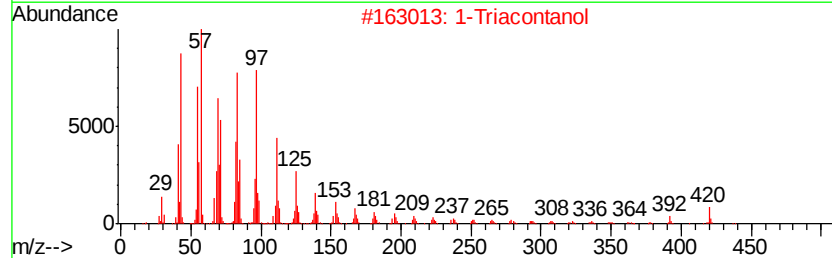
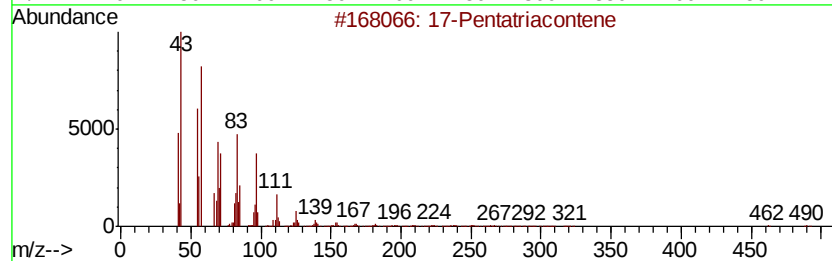
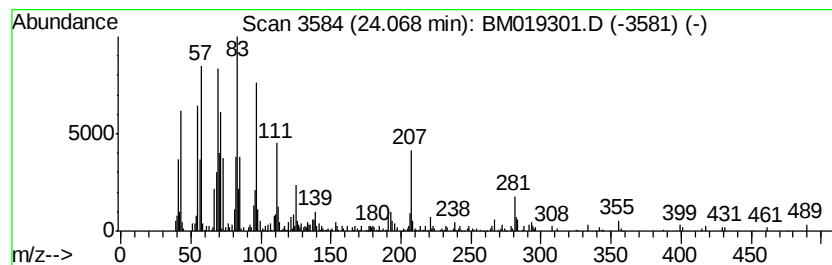
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM031119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 14 17-Pentatriacontene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.07	3.25 ng	439894	Perylene-d12	23.47

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	17-Pentatriacontene	491	C35H70	006971-40-0	83
2		1-Triacontanol	438	C30H62O	000593-50-0	60
3		Isoheptadecanol	256	C17H36O	057289-07-3	58
4		Thieno[2,3-d]-1,3-thiaselenol-2-...	238	C5H2S3Se	081803-09-0	56
5		Octadecane, 1-(ethenyloxy)-	296	C20H40O	000930-02-9	52



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032119\
Data File : BM019301.D
Acq On : 21 Mar 2019 21:33
Operator : JU/SJ
Sample : K2004-04
Misc :
ALS Vial : 13 Sample Multiplier: 1

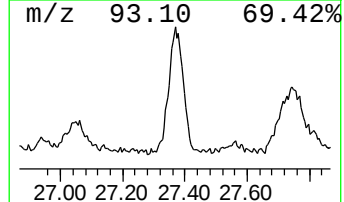
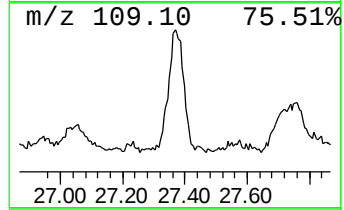
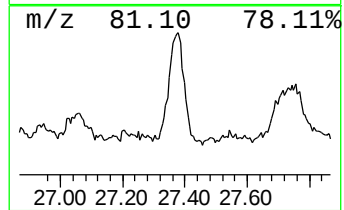
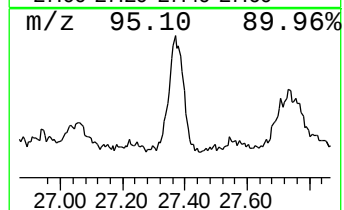
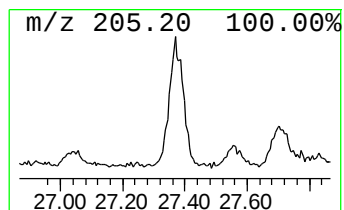
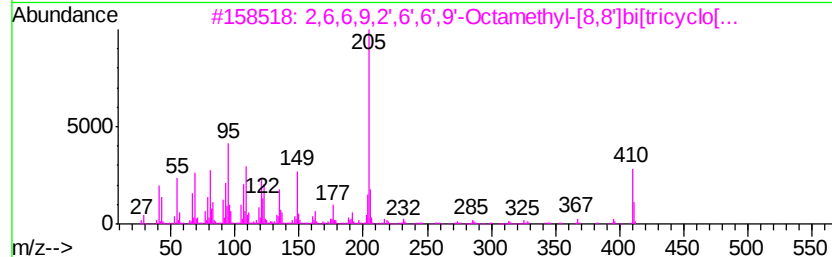
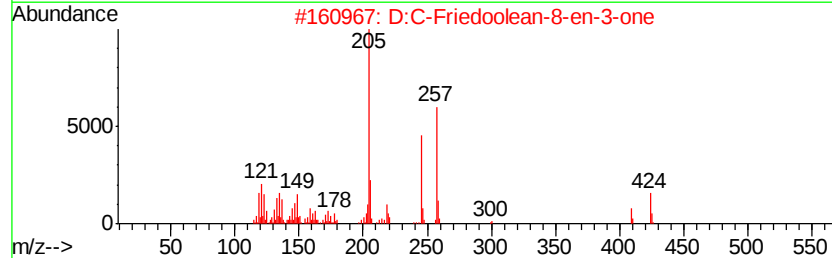
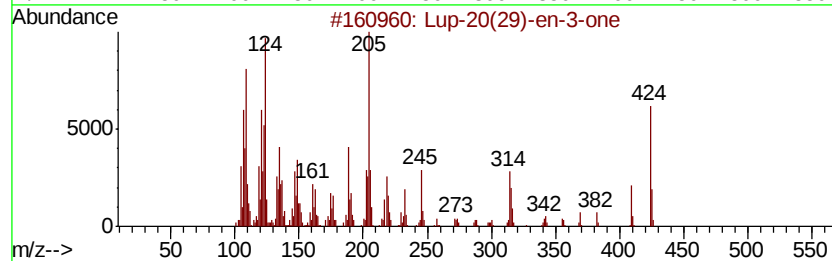
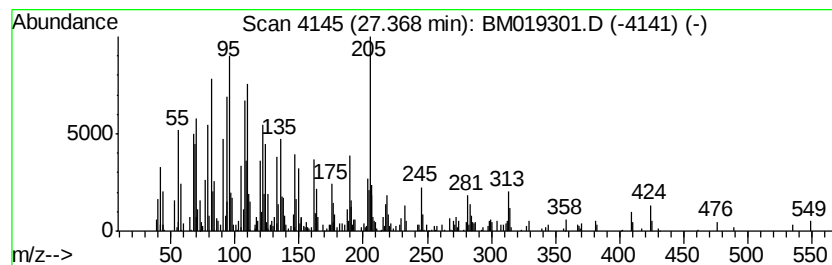
Instrument :
BNA_M
ClientSampleId :
SB-19(13-14)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM031119.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Lup-20(29)-en-3-one Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
27.37	9.04 ng	1224950	Perylene-d12	23.47	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Lup-20(29)-en-3-one	424	C30H48O	001617-70-5	81
2	D:C-Friedoolean-8-en-3-one	424	C30H48O	022611-26-3	70
3	2,6,6,9,2',6',6',9'-Octamethyl-[...]	410	C30H50	1000189-48-2	53
4	9,19-Cyclolanost-23-ene-3,25-dio...	442	C30H50O2	054482-57-4	45
5	Bicyclo[2.2.1]heptane, 2,2,3-tri...	138	C10H18	000473-19-8	41



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM032119\
 Data File : BM019301.D
 Acq On : 21 Mar 2019 21:33
 Operator : JU/SJ
 Sample : K2004-04
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SB-19(13-14)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\8270-BM031119.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
unknown7.16	7.16	112.5	ng	7498620	1	7.63	1333110	20.0
Phenol, 3-ethyl-	9.87	9.6	ng	908016	2	10.41	1887090	20.0
1,3-Cyclopentadie...	11.25	2.2	ng	204324	2	10.41	1887090	20.0
n-Hexadecanoic acid	17.94	8.6	ng	1374620	4	17.04	3206730	20.0
dl-3-Indolelactic...	18.42	2.1	ng	330581	4	17.04	3206730	20.0
1H-Indole-3-metha...	18.61	2.0	ng	321243	4	17.04	3206730	20.0
Octadecanoic acid	19.22	3.8	ng	479271	5	21.24	2539970	20.0
Cyclotetradecane	20.95	13.9	ng	1763680	5	21.24	2539970	20.0
unknown21.76	21.76	2.5	ng	314961	5	21.24	2539970	20.0
Heneicosane	21.81	2.1	ng	272036	5	21.24	2539970	20.0
1-Eicosanol	21.84	5.7	ng	720924	5	21.24	2539970	20.0
Tetradecane	22.77	2.8	ng	376198	6	23.47	2711100	20.0
17-Pentatriacontene	24.07	3.3	ng	439894	6	23.47	2711100	20.0
Lup-20(29)-en-3-one	27.37	9.0	ng	1224950	6	23.47	2711100	20.0