

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121618\
 Data File : BM018224.D
 Acq On : 16 Dec 2018 13:37
 Operator : JU/SJ
 Sample : J6377-03
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 P001-SD014-0006-01

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-BM121518.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	4.687	302	306	313	rBV	101548	138115	1.09%	0.212%
2	4.869	332	337	346	rBV	246262	336849	2.67%	0.516%
3	5.134	375	382	386	rBV	2953410	4126702	32.66%	6.325%
4	5.187	386	391	403	rVB	9068621	12636665	100.00%	19.369%
5	6.592	625	630	632	rBV	118873	175159	1.39%	0.268%
6	6.628	632	636	642	rVV	655477	928331	7.35%	1.423%
7	6.698	642	648	664	rVB	2701892	4695462	37.16%	7.197%
8	7.040	698	706	715	rBV	3015908	4567463	36.14%	7.001%
9	7.145	718	724	731	rVB	301779	453285	3.59%	0.695%
10	7.498	778	784	787	rBV	649281	1017974	8.06%	1.560%
11	7.534	787	790	796	rVV	432551	623199	4.93%	0.955%
12	7.604	796	802	810	rVB2	90096	184383	1.46%	0.283%
13	7.810	830	837	850	rVB	2105683	3255336	25.76%	4.990%
14	8.657	973	981	999	rBV	1601369	2695422	21.33%	4.131%
15	9.033	1039	1045	1052	rVB	127430	190254	1.51%	0.292%
16	10.263	1244	1254	1260	rBV	800070	1383607	10.95%	2.121%
17	10.316	1260	1263	1270	rVB	104396	161765	1.28%	0.248%
18	12.763	1670	1679	1688	rBV	3790676	5456975	43.18%	8.364%
19	13.627	1820	1826	1832	rBV	205164	332734	2.63%	0.510%
20	14.157	1909	1916	1924	rBV2	1178162	1794502	14.20%	2.751%
21	15.662	2164	2172	2180	rBV	2939544	4201793	33.25%	6.440%
22	15.910	2206	2214	2219	rBV3	94606	204963	1.62%	0.314%
23	16.915	2379	2385	2391	rBV2	1250670	1816413	14.37%	2.784%
24	16.986	2395	2397	2402	rVB	106666	130969	1.04%	0.201%
25	17.109	2413	2418	2423	rBV	92478	139481	1.10%	0.214%
26	17.833	2537	2541	2550	rVB4	231157	364734	2.89%	0.559%
27	19.127	2757	2761	2767	rBV5	114608	199379	1.58%	0.306%
28	19.574	2831	2837	2844	rVB	4318543	5330970	42.19%	8.171%
29	19.850	2881	2884	2888	rBV3	134502	182541	1.44%	0.280%
30	20.133	2928	2932	2937	rVB	165265	205659	1.63%	0.315%
31	20.862	3052	3056	3061	rBV2	339861	439076	3.47%	0.673%
32	20.974	3072	3075	3079	rBV2	88872	132615	1.05%	0.203%
33	21.139	3098	3103	3112	rBV	1310017	1895430	15.00%	2.905%
34	21.297	3127	3130	3138	rVB	167811	198846	1.57%	0.305%

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Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

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Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	21.721	3199	3202	3205	rBV	334773	411702	3.26%	0.631%
36	21.750	3205	3207	3213	rVB3	289491	390846	3.09%	0.599%
37	22.168	3274	3278	3282	rVB	252658	322168	2.55%	0.494%
38	22.650	3355	3360	3364	rBV	684376	969422	7.67%	1.486%
39	22.697	3366	3368	3378	rVB7	132583	234131	1.85%	0.359%
40	23.186	3447	3451	3457	rVB	267535	371393	2.94%	0.569%
41	23.303	3465	3471	3478	rVB	864113	1554588	12.30%	2.383%
42	23.791	3549	3554	3559	rVB2	221262	389514	3.08%	0.597%

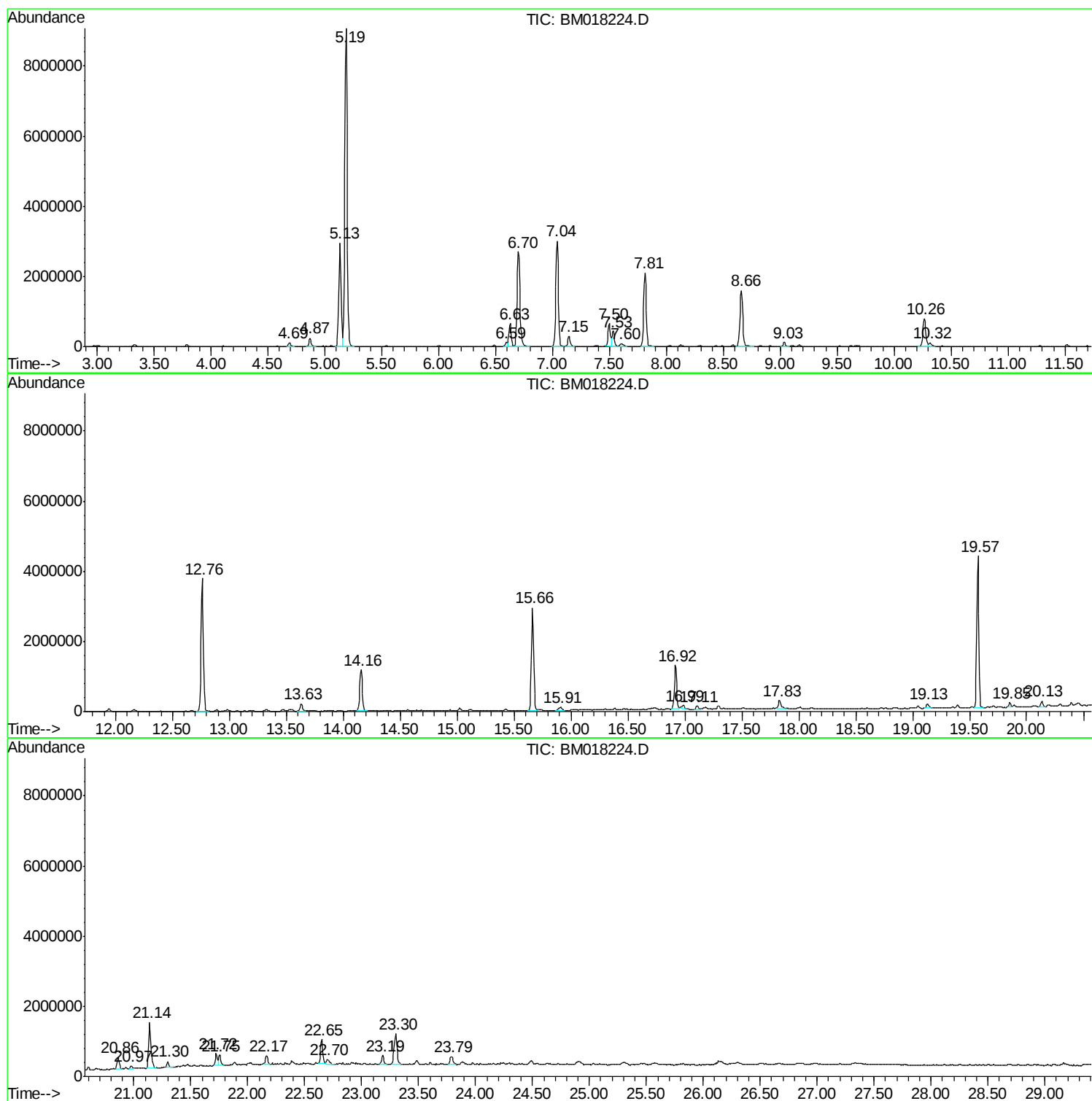
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM121518.M
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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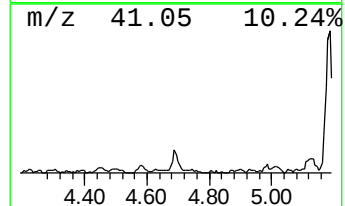
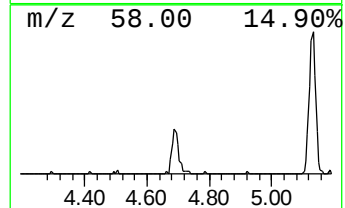
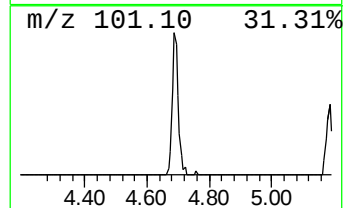
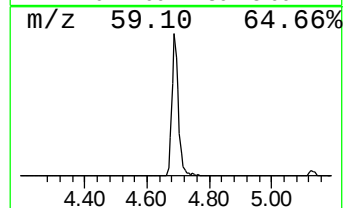
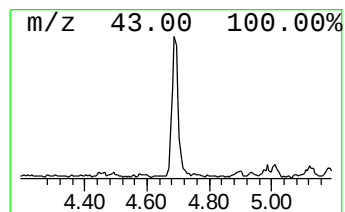
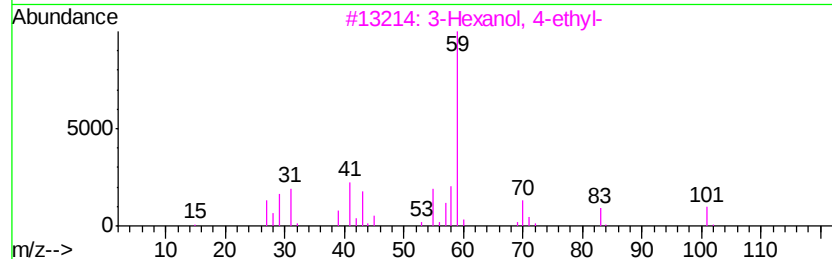
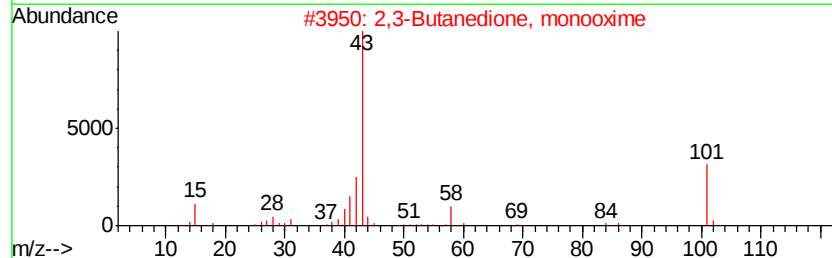
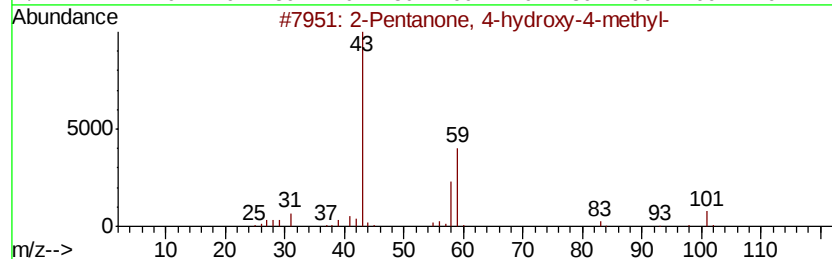
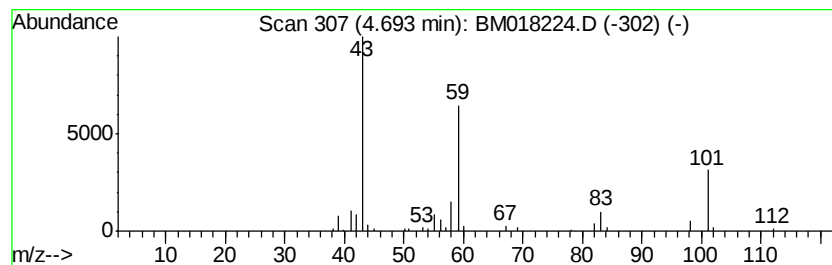
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM121518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.69	2.71 ng	138115	1,4-Dichlorobenzene-d4	7.50

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	35
3		3-Hexanol, 4-ethyl-	130	C8H18O	019780-44-0	33
4		5-Ethyl-3-methylhept-1-en-4-ol	156	C10H20O	286424-80-4	28
5		3-Octanol	130	C8H18O	000589-98-0	25



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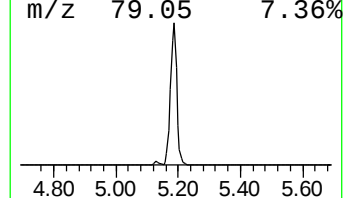
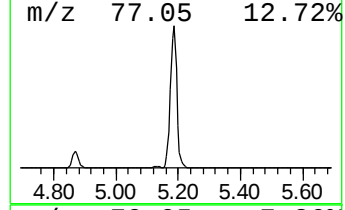
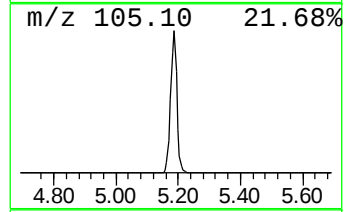
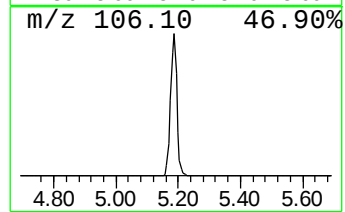
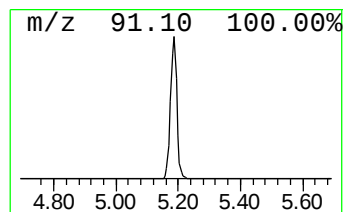
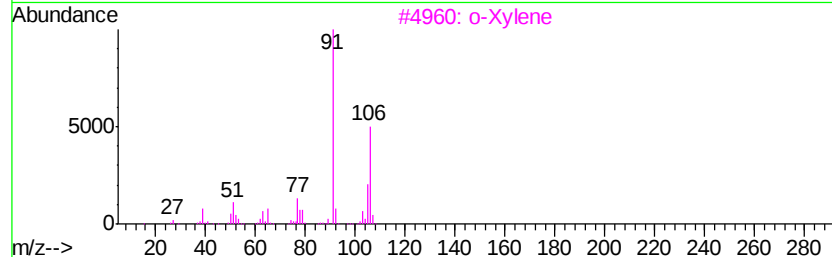
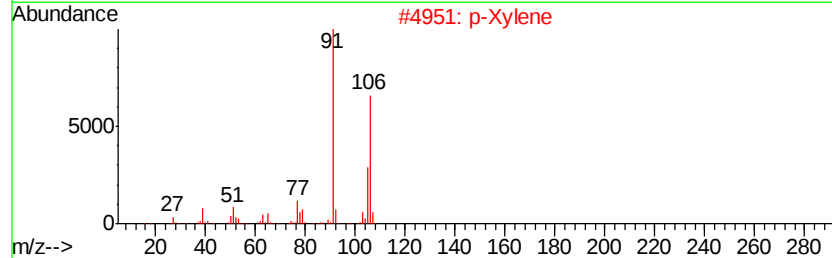
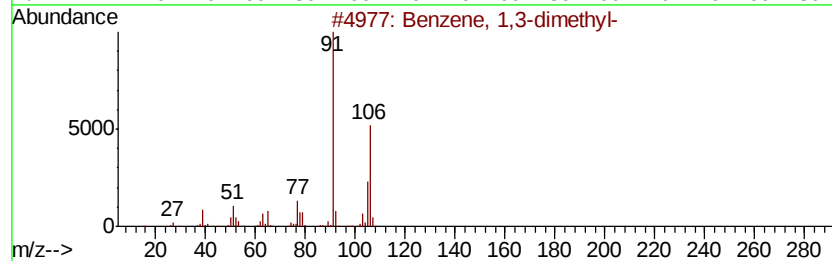
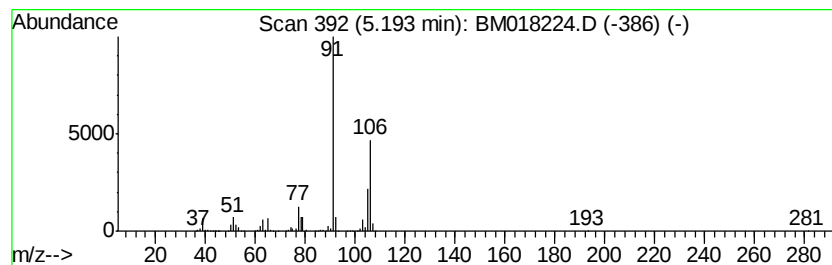
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 3 Benzene, 1,3-dimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.19	248.27 ng	12636700	1,4-Dichlorobenzene-d4	7.50

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	97
2		p-Xylene	106	C8H10	000106-42-3	95
3		o-Xylene	106	C8H10	000095-47-6	93
4		1,3-Cyclopentadiene, 5-(1-methyl...	106	C8H10	002175-91-9	90
5		Ethylbenzene	106	C8H10	000100-41-4	83



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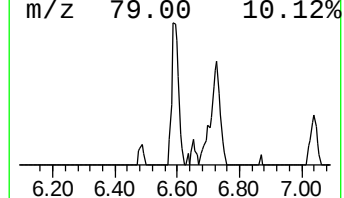
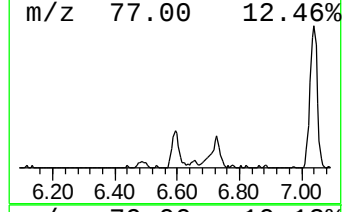
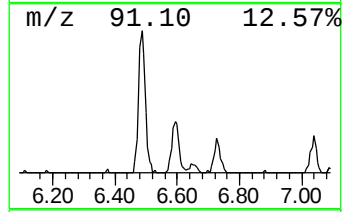
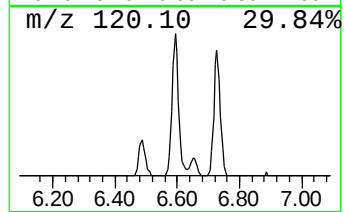
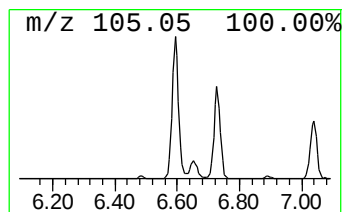
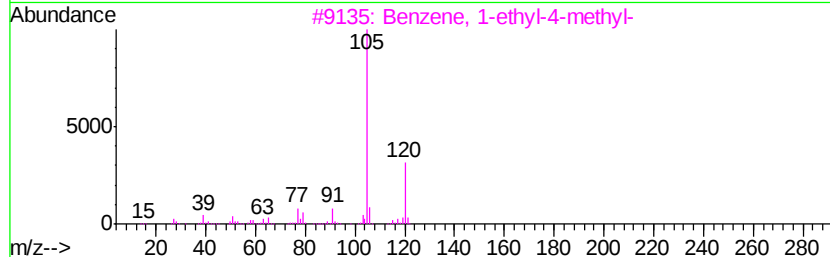
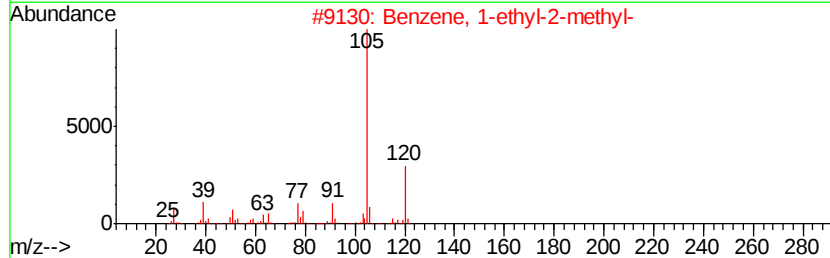
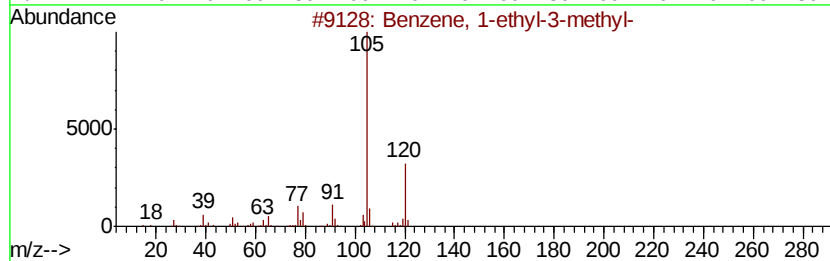
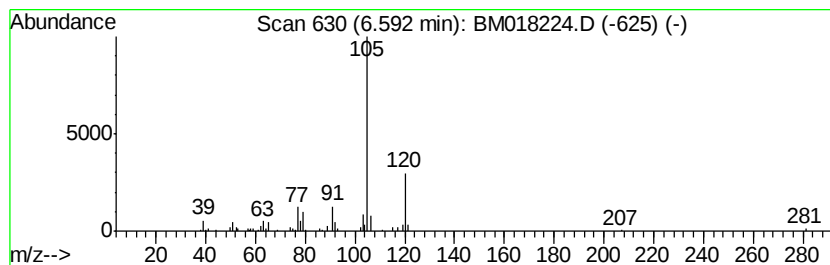
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Benzene, 1-ethyl-3-methyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.59	3.44 ng	175159	1,4-Dichlorobenzene-d4	7.50

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	97
2		Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
3		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
4		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
5		Benzene, 1,3,5-trimethyl-	120	C9H12	000108-67-8	91



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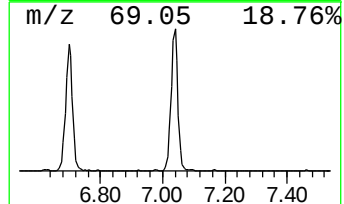
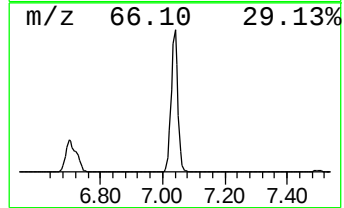
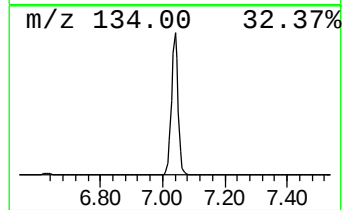
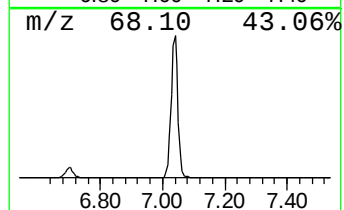
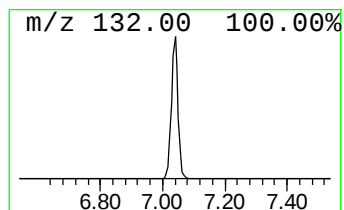
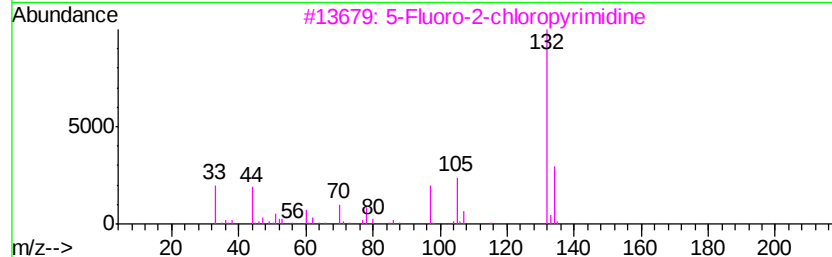
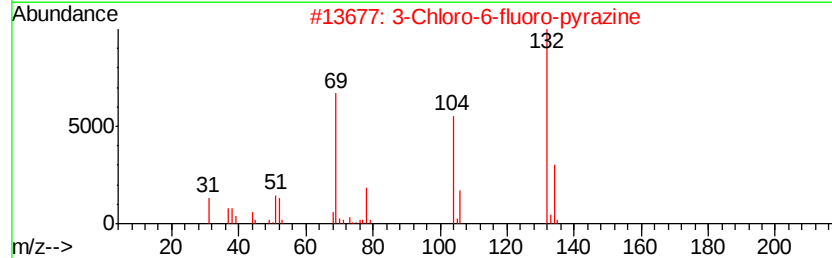
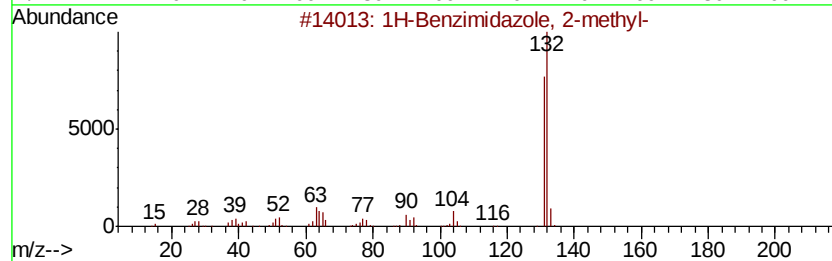
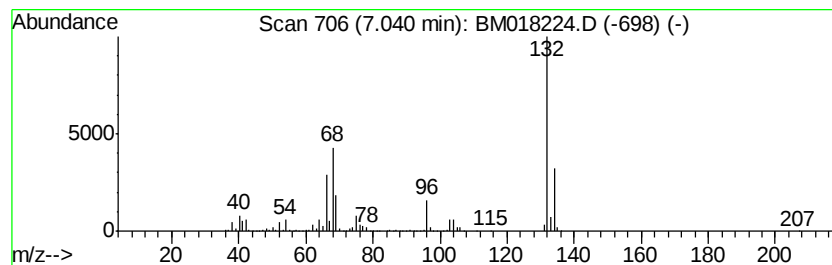
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 unknown7.04 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.04	89.74 ng	4567460	1,4-Dichlorobenzene-d4	7.50

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		(5-METHYL-2-PYRIDYL)ACETONITRILE	132	C8H8N2	1000241-93-9	10
5		Xenon	132	Xe	007440-63-3	9



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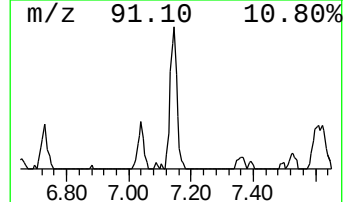
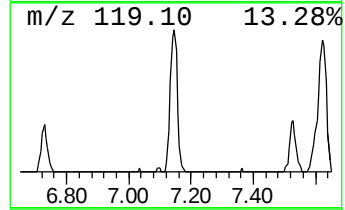
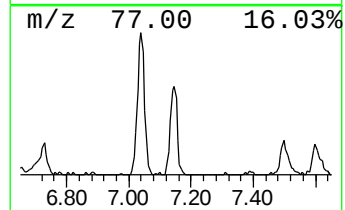
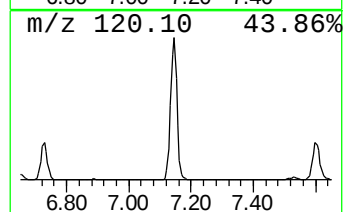
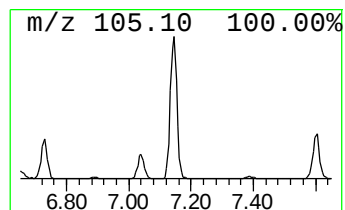
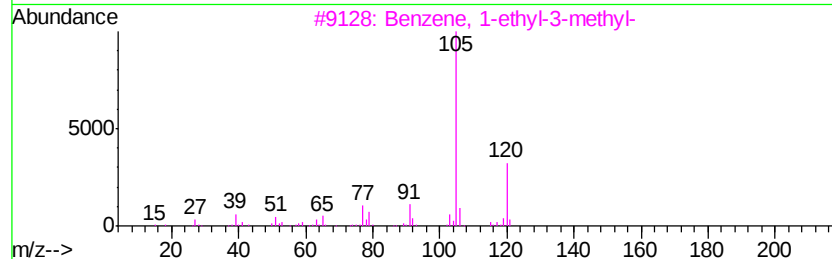
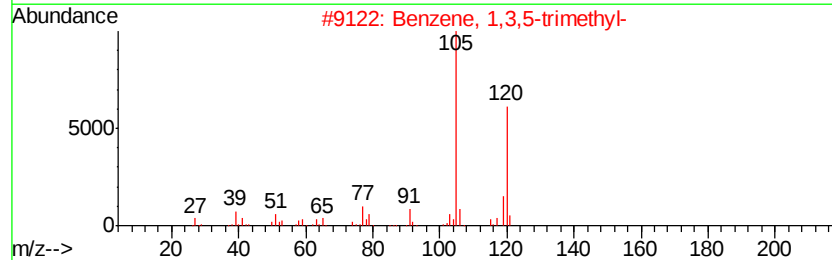
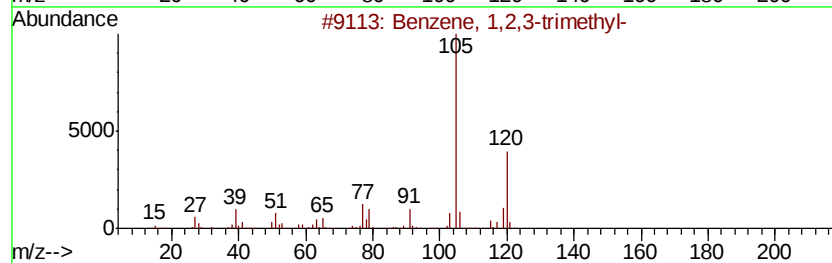
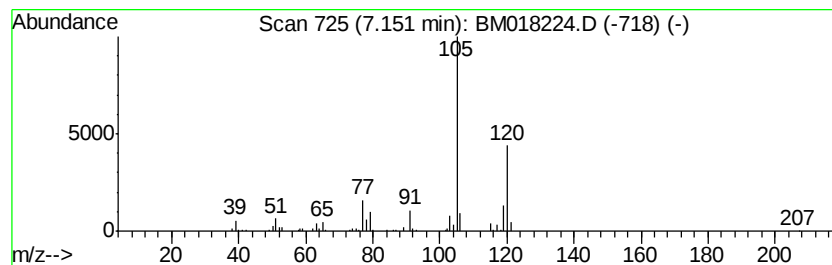
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Benzene, 1,2,3-trimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.15	8.91 ng	453285	1,4-Dichlorobenzene-d4	7.50

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
2		Benzene, 1,3,5-trimethyl-	120	C9H12	000108-67-8	94
3		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
4		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121618\
Data File : BM018224.D
Acq On : 16 Dec 2018 13:37
Operator : JU/SJ
Sample : J6377-03
Misc :
ALS Vial : 6 Sample Multiplier: 1

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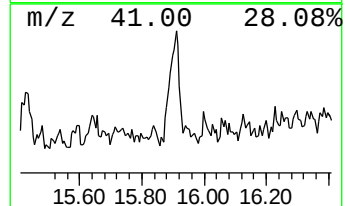
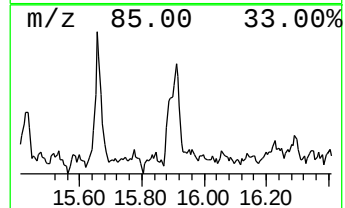
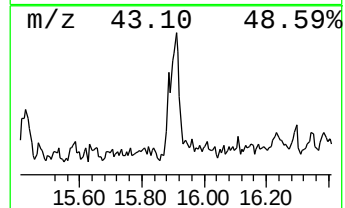
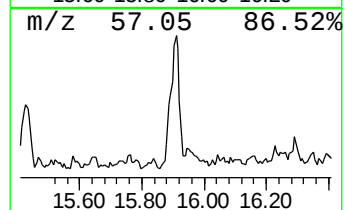
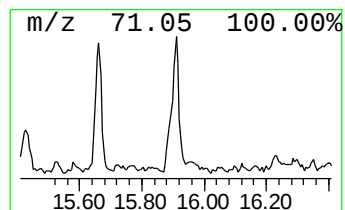
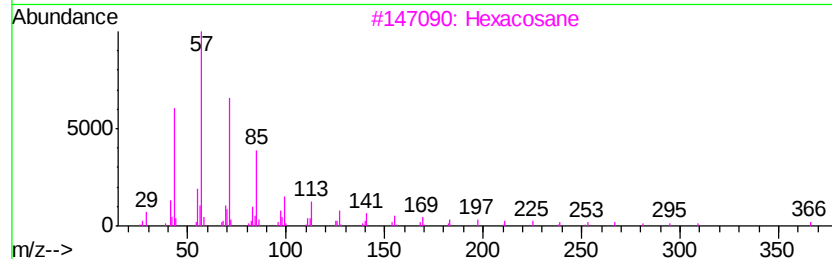
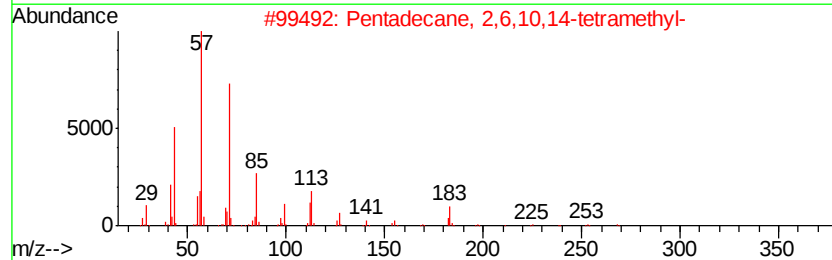
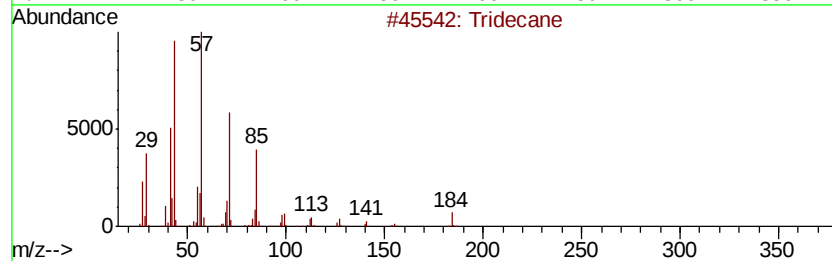
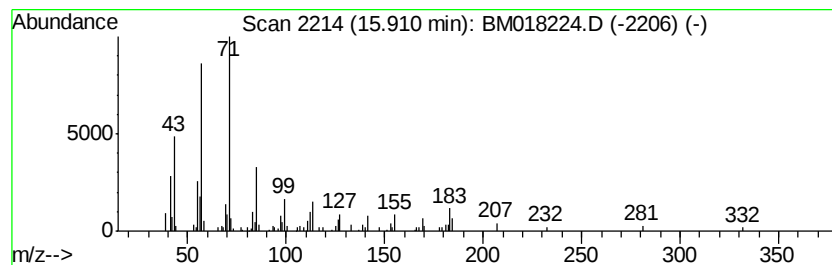
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Tridecane Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.91	2.26 ng	204963	Phenanthrene-d10	16.92

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tridecane	184	C13H28	000629-50-5	72
2		Pentadecane, 2,6,10,14-tetramethyl-	268	C19H40	001921-70-6	58
3		Hexacosane	366	C26H54	000630-01-3	53
4		Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	52
5		Hentriacontane	437	C31H64	000630-04-6	49



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121618\
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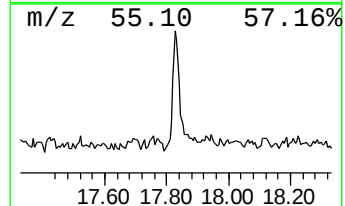
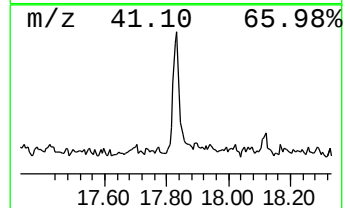
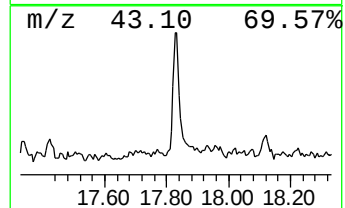
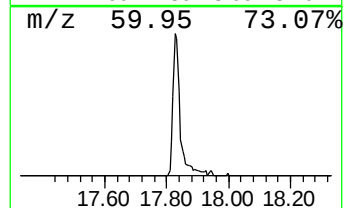
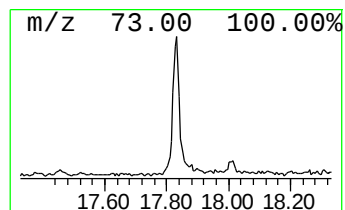
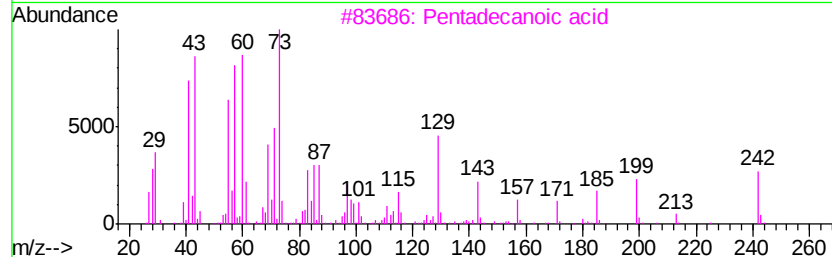
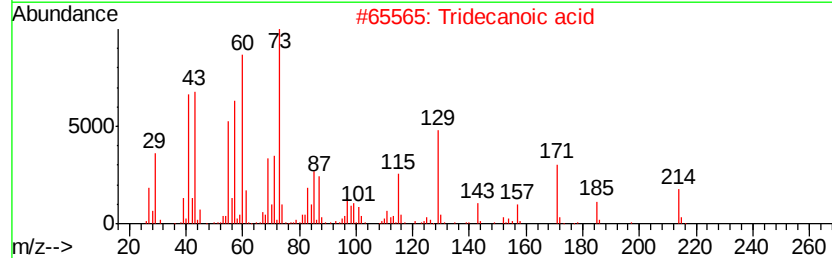
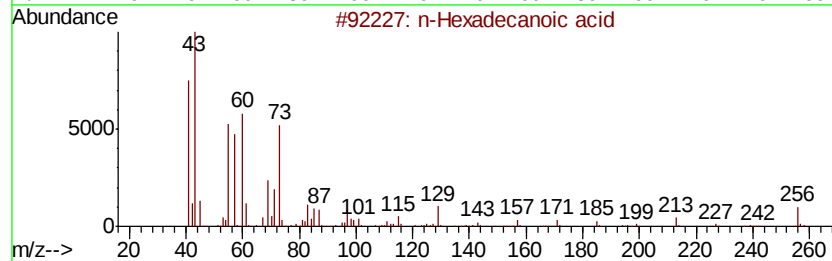
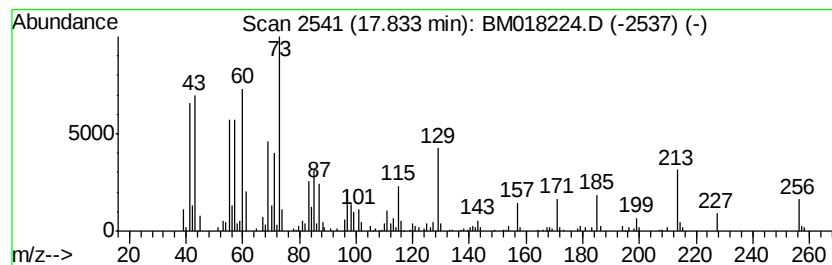
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 n-Hexadecanoic acid Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.	
17.83	4.02 ng	364734	Phenanthrene-d10	16.92	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
2	Tridecanoic acid	214	C13H26O2	000638-53-9	90
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	83
4	n-Decanoic acid	172	C10H20O2	000334-48-5	62
5	Estra-1,3,5(10)-trien-17.beta.-ol	256	C18H24O	002529-64-8	30



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121618\
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Sample : J6377-03
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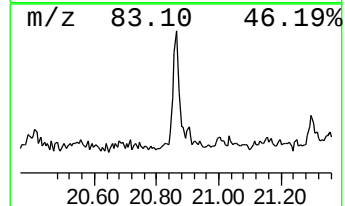
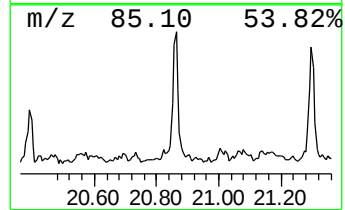
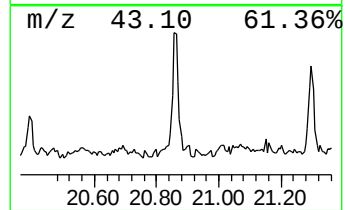
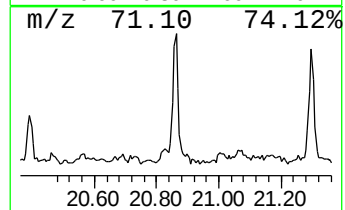
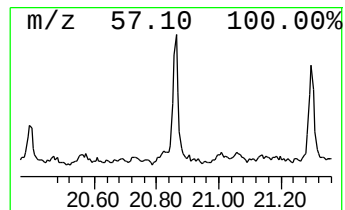
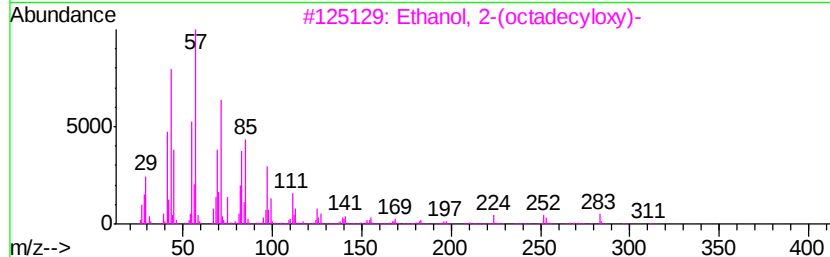
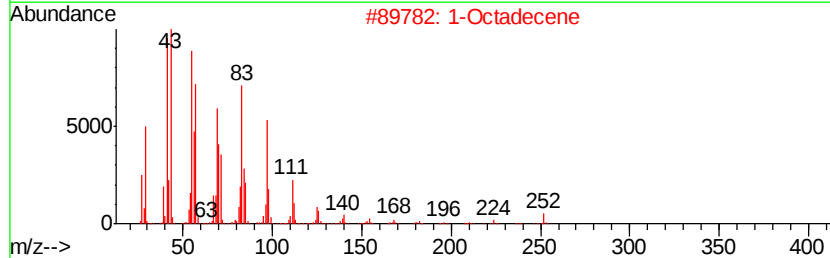
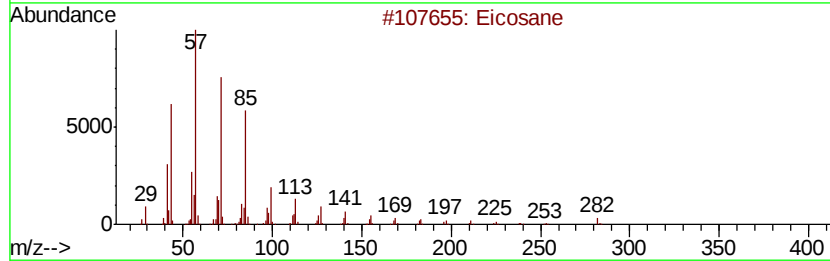
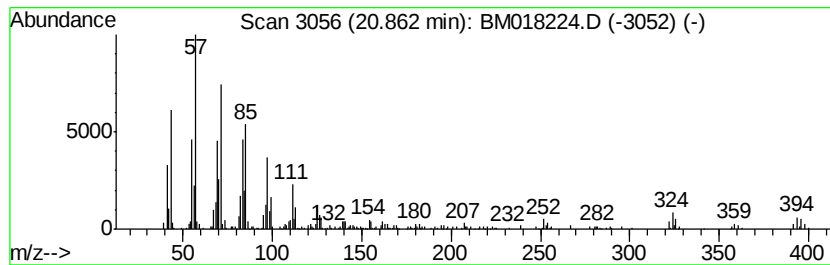
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Eicosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.86	4.63 ng	439076	Chrysene-d12	21.14
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Eicosane	282 C20H42	000112-95-8	90
2	1-Octadecene	252 C18H36	000112-88-9	89
3	Ethanol, 2-(octadecyloxy)-	314 C20H42O2	002136-72-3	83
4	Octadecane, 1-[2-(hexadecyloxy)e...	539 C36H74O2	017367-10-1	80
5	Octadecane	254 C18H38	000593-45-3	70



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121618\
Data File : BM018224.D
Acq On : 16 Dec 2018 13:37
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Misc :
ALS Vial : 6 Sample Multiplier: 1

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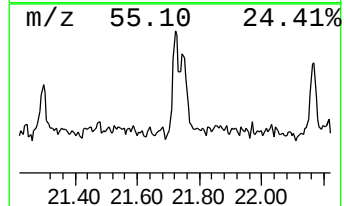
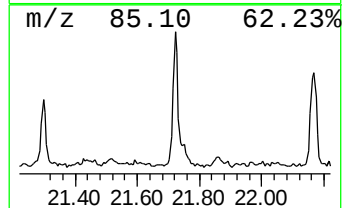
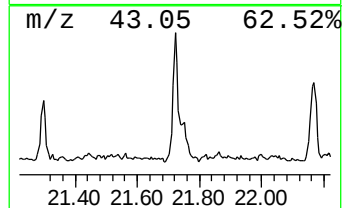
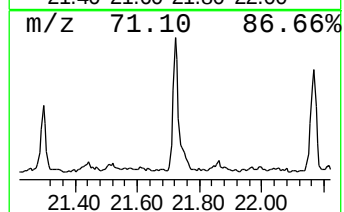
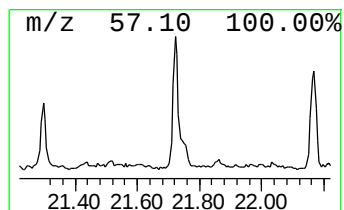
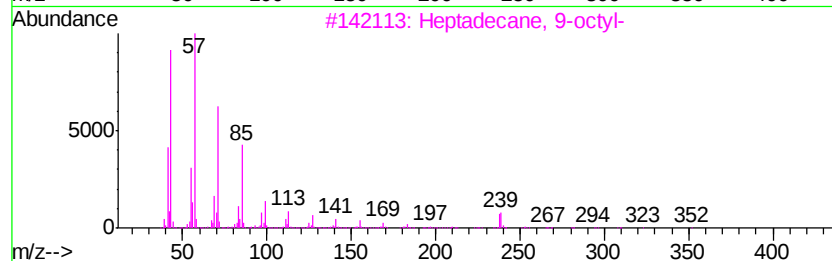
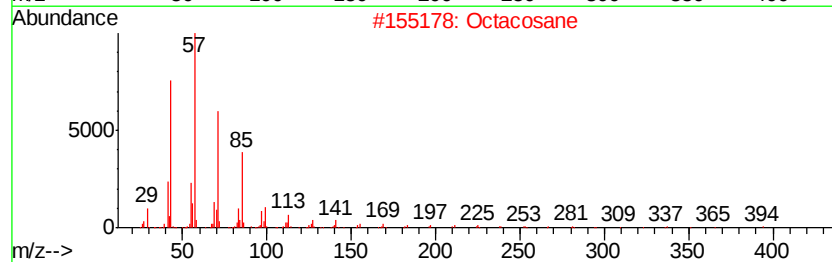
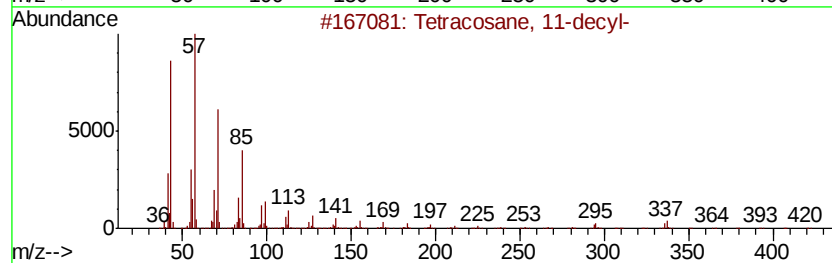
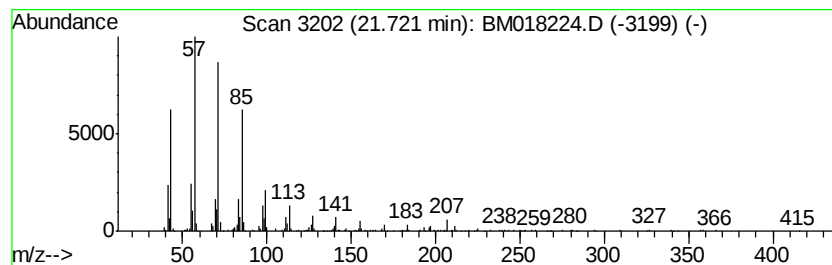
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Tetracosane, 11-decyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.72	4.34 ng	411702	Chrysene-d12	21.14

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetracosane, 11-decyl-	479	C34H70	055429-84-0	91
2			Octacosane	394	C28H58	000630-02-4	91
3			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
4			Heneicosane	296	C21H44	000629-94-7	90
5			Tetratriacontane	479	C34H70	014167-59-0	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121618\
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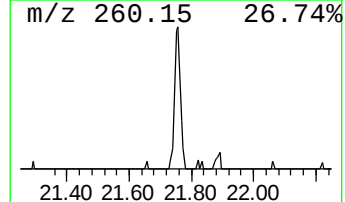
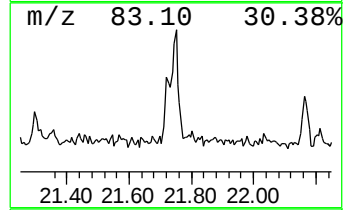
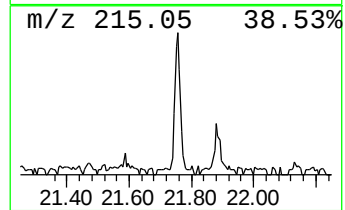
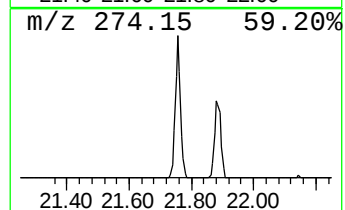
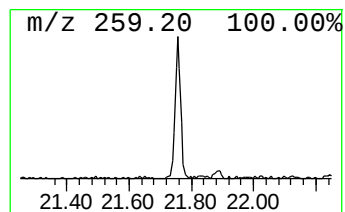
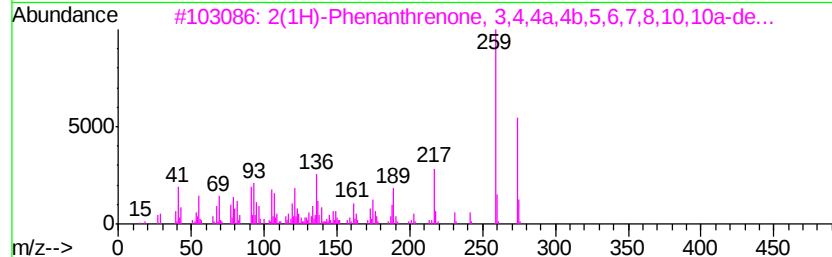
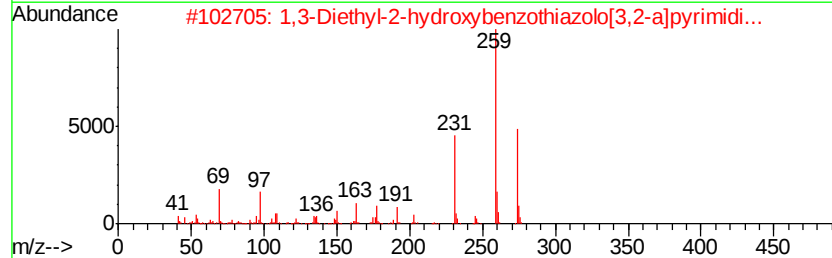
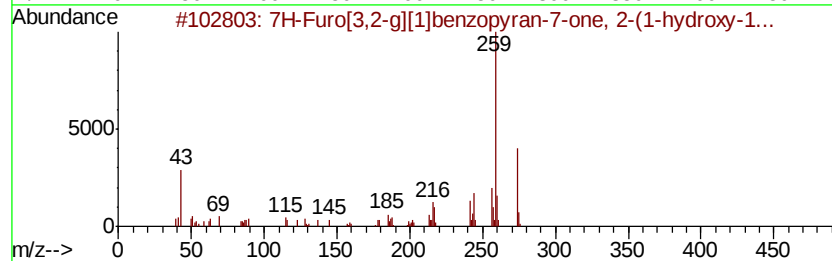
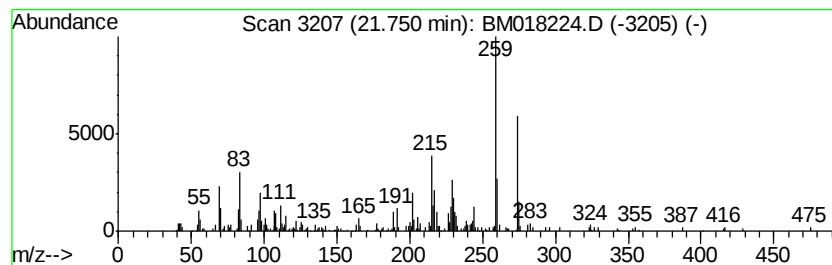
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 7H-Furo[3,2-g][1]benzopyran... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.		
21.75	4.12 ng	390846	Chrysene-d12	21.14		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		7H-Furo[3,2-g][1]benzopyran-7-on...	274	C15H14O5	054087-32-0	53
2		1,3-Diethyl-2-hydroxybenzothiazo...	274	C14H14N2O2S	1000227-15-3	50
3		2(1H)-Phenanthrenone, 3,4,4a,4b,...	274	C19H30O	007715-44-8	47
4		Androstan-6-one, (5.alpha.)-	274	C19H30O	003676-06-0	38
5		5.beta.-Androstan-3.beta.,17.bet...	292	C19H32O2	006038-31-9	38



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121618\
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Acq On : 16 Dec 2018 13:37
Operator : JU/SJ
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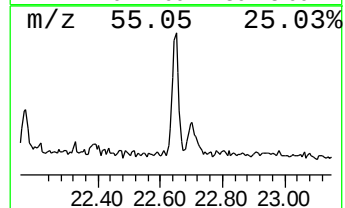
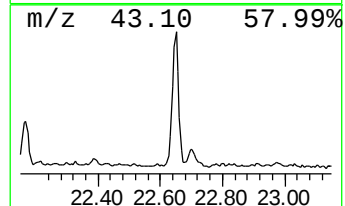
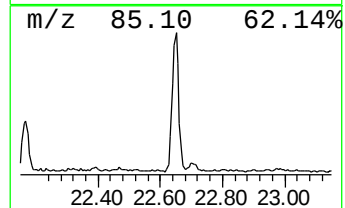
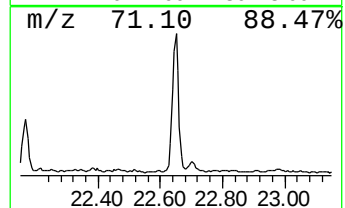
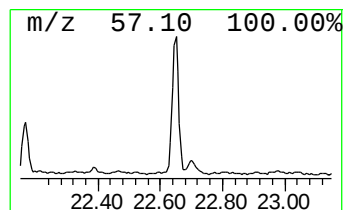
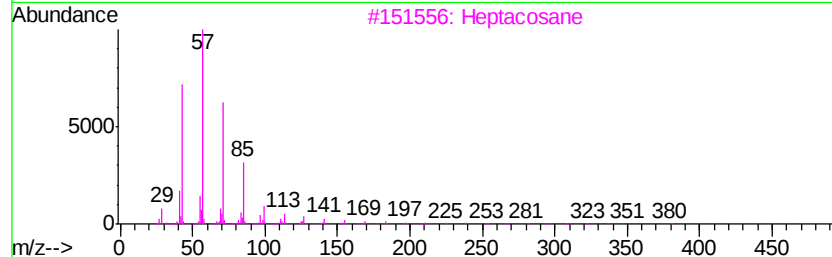
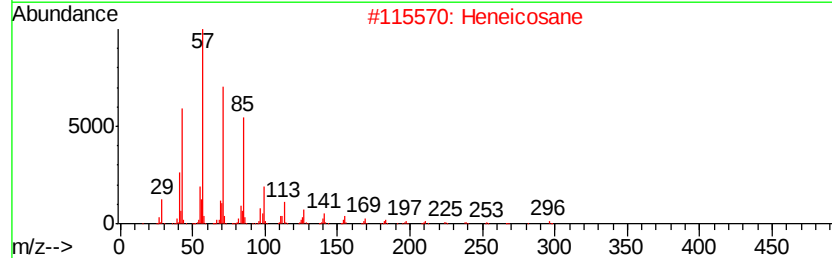
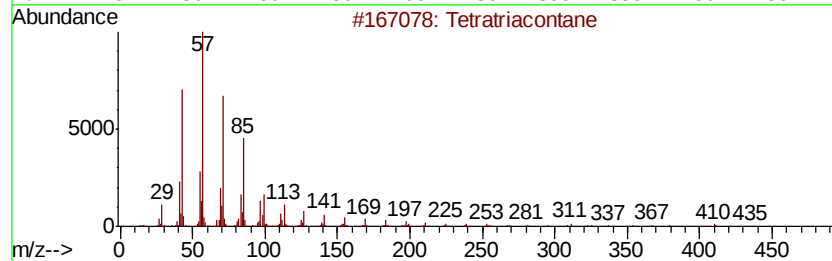
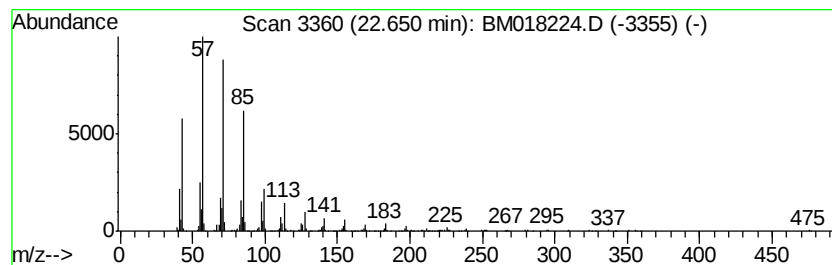
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 17 Tetratriacontane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.65	12.47 ng	969422	Perylene-d12	23.30

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetratriacontane	479	C34H70	014167-59-0	91
2		Heneicosane	296	C21H44	000629-94-7	91
3		Heptacosane	380	C27H56	000593-49-7	91
4		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
5		Triacontane	422	C30H62	000638-68-6	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121618\
Data File : BM018224.D
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Sample : J6377-03
Misc :
ALS Vial : 6 Sample Multiplier: 1

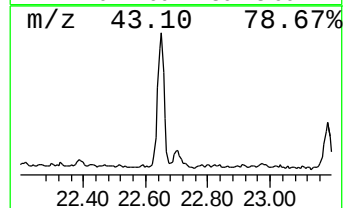
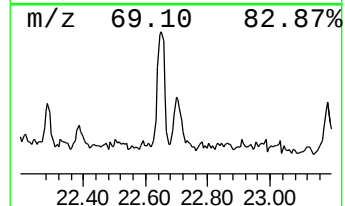
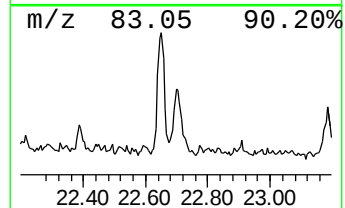
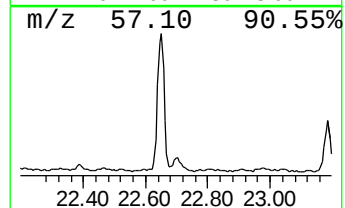
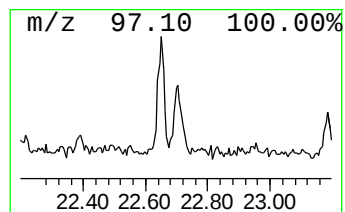
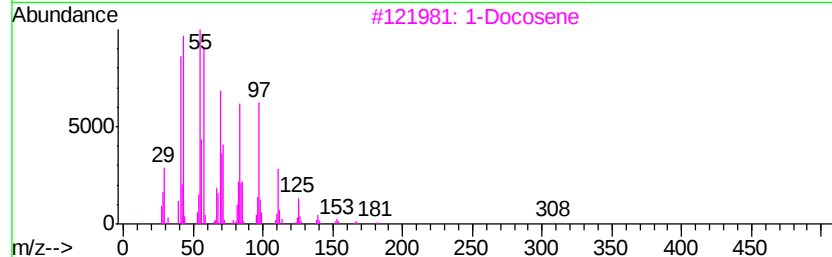
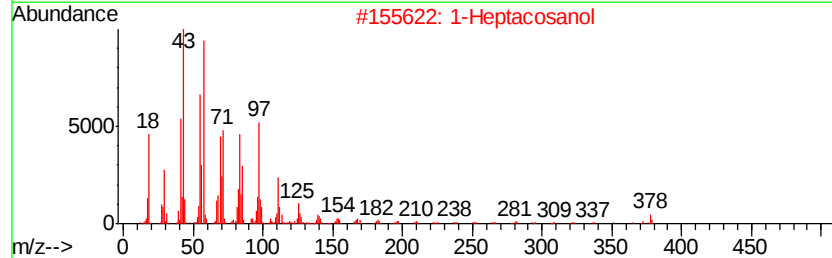
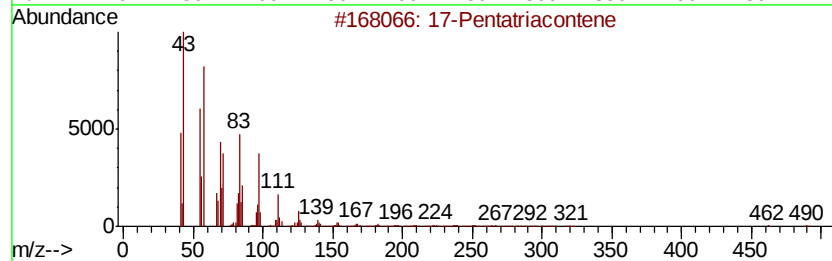
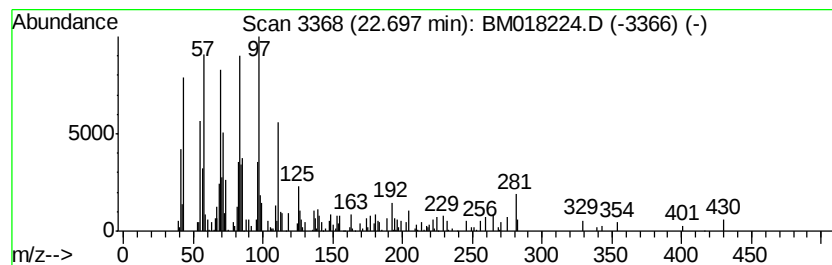
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BNA_M
ClientSampleId :
P001-SD014-0006-01

Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM121518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 17-Pentatriacontene Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.	
22.70	3.01 ng	234131	Perylene-d12	23.30	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	17-Pentatriacontene	491	C35H70	006971-40-0	86
2	1-Heptacosanol	396	C27H56O	002004-39-9	53
3	1-Docosene	308	C22H44	001599-67-3	49
4	Heptafluorobutanoic acid, heptadecanoic acid	452	C21H35F7O2	1000282-97-3	49
5	1H-Imidazole, 1-(cyclohexylacetyl)-	192	C11H16N2O	093383-50-7	43



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121618\
Data File : BM018224.D
Acq On : 16 Dec 2018 13:37
Operator : JU/SJ
Sample : J6377-03
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
P001-SD014-0006-01

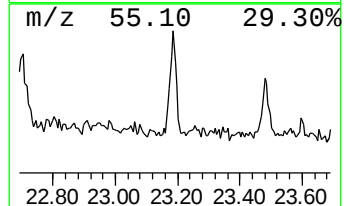
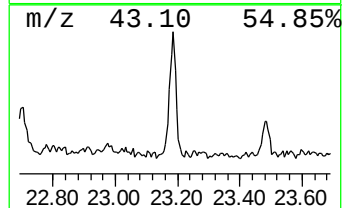
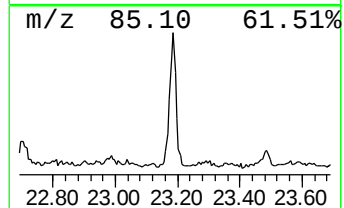
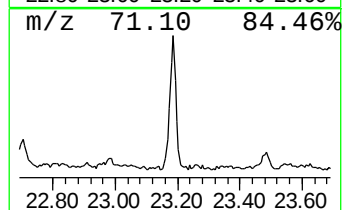
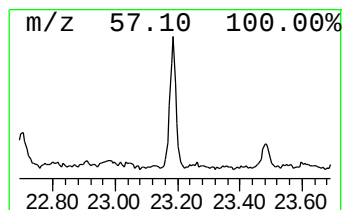
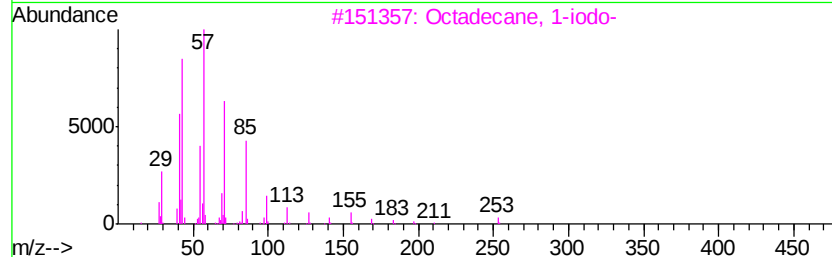
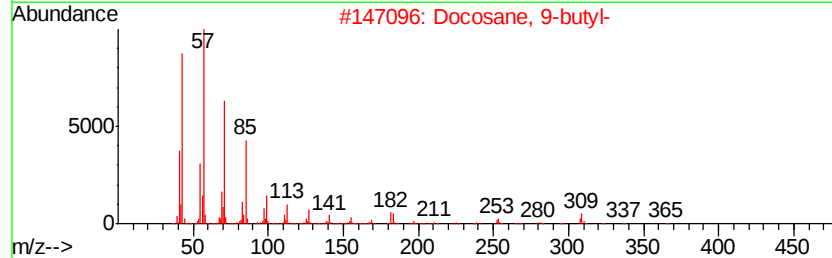
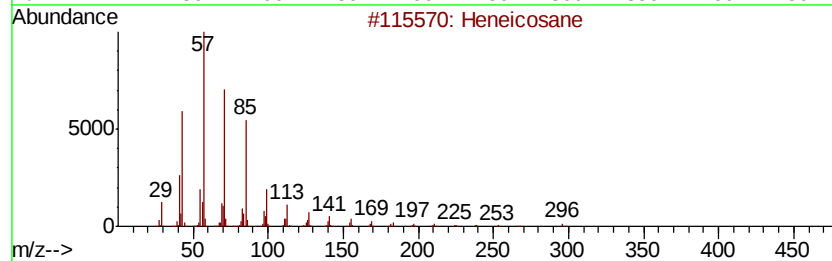
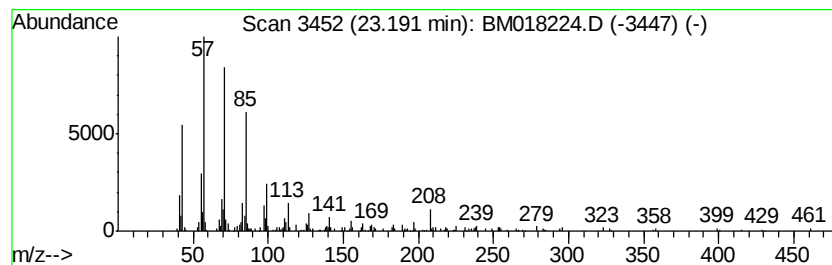
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 Heneicosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.19	4.78 ng	371393	Perylene-d12	23.30

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	91
2		Docosane, 9-butyl-	366	C26H54	055282-14-9	90
3		Octadecane, 1-iodo-	380	C18H37I	000629-93-6	87
4		Nonacosane	408	C29H60	000630-03-5	87
5		Hexadecane	226	C16H34	000544-76-3	83



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121618\
Data File : BM018224.D
Acq On : 16 Dec 2018 13:37
Operator : JU/SJ
Sample : J6377-03
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
P001-SD014-0006-01

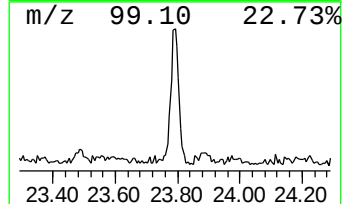
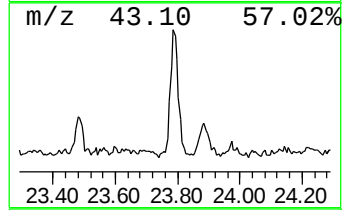
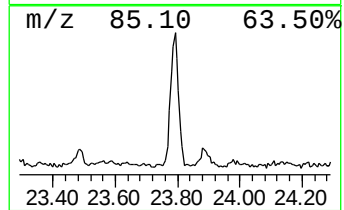
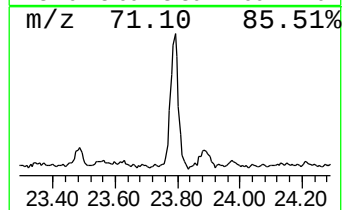
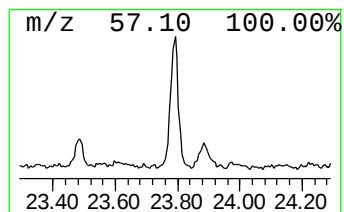
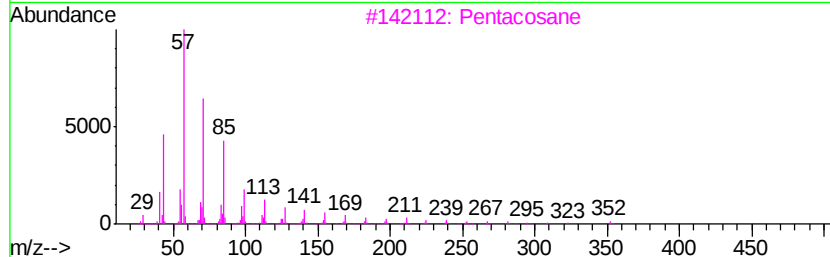
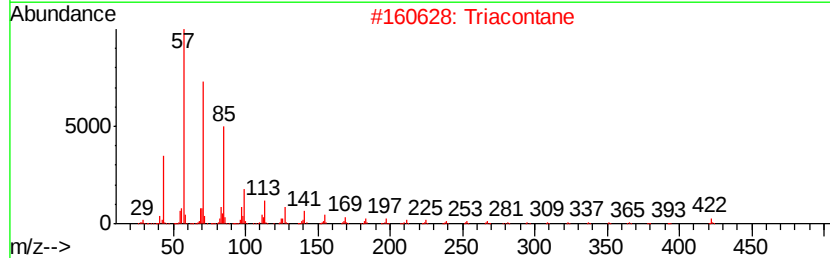
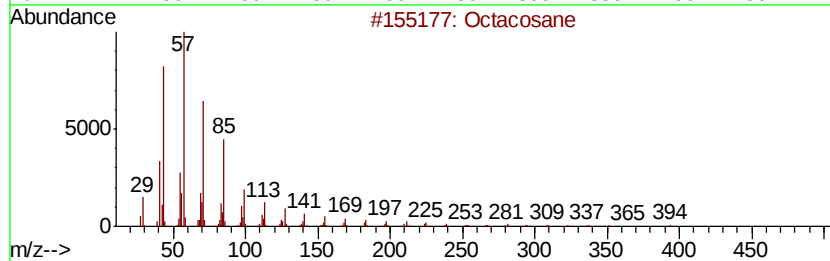
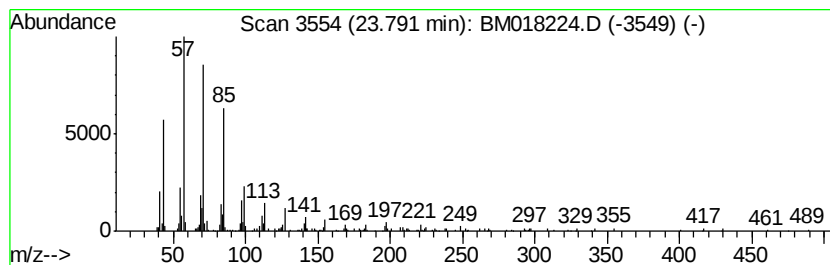
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 20 Octacosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.79	5.01 ng	389514	Perylene-d12	23.30

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octacosane	394	C28H58	000630-02-4	91
2		Triacontane	422	C30H62	000638-68-6	91
3		Pentacosane	352	C25H52	000629-99-2	90
4		Heneicosane	296	C21H44	000629-94-7	90
5		Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM121618\
 Data File : BM018224.D
 Acq On : 16 Dec 2018 13:37
 Operator : JU/SJ
 Sample : J6377-03
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 P001-SD014-0006-01

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\8270-BM121518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.69	2.7	ng	138115	1	7.50	1017970	20.0
Benzene, 1,3-dime...	5.19	248.3	ng	12636700	1	7.50	1017970	20.0
Benzene, 1-ethyl-...	6.59	3.4	ng	175159	1	7.50	1017970	20.0
unknown7.04	7.04	89.7	ng	4567460	1	7.50	1017970	20.0
Benzene, 1,2,3-tr...	7.15	8.9	ng	453285	1	7.50	1017970	20.0
Tridecane	15.91	2.3	ng	204963	4	16.92	1816410	20.0
n-Hexadecanoic acid	17.83	4.0	ng	364734	4	16.92	1816410	20.0
Eicosane	20.86	4.6	ng	439076	5	21.14	1895430	20.0
Tetracosane, 11-d...	21.72	4.3	ng	411702	5	21.14	1895430	20.0
7H-Furo[3,2-g][1]...	21.75	4.1	ng	390846	5	21.14	1895430	20.0
Tetratriacontane	22.65	12.5	ng	969422	6	23.30	1554590	20.0
17-Pentatriacontene	22.70	3.0	ng	234131	6	23.30	1554590	20.0
Heneicosane	23.19	4.8	ng	371393	6	23.30	1554590	20.0
Octacosane	23.79	5.0	ng	389514	6	23.30	1554590	20.0