

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM021819\
 Data File : BM018843.D
 Acq On : 18 Feb 2019 18:00
 Operator : JU/SJ
 Sample : K1518-19
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 TP-MH-C05

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-BM021119.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	4.857	297	301	312	rVB	335833	486694	2.49%	0.336%
2	5.299	369	376	391	rBV	7810941	11062057	56.66%	7.628%
3	6.887	638	646	663	rBV	7398839	12657771	64.84%	8.729%
4	7.240	698	706	722	rBV	7900216	12238252	62.69%	8.439%
5	7.704	778	785	793	rBV	1323947	2060323	10.55%	1.421%
6	8.022	831	839	850	rBV	5349696	8447221	43.27%	5.825%
7	8.875	975	984	996	rBV	4359241	7133344	36.54%	4.919%
8	10.492	1247	1259	1266	rBV	1712989	2847793	14.59%	1.964%
9	12.969	1673	1680	1696	rBV	10157056	14752869	75.57%	10.174%
10	13.816	1817	1824	1836	rBV	407601	612580	3.14%	0.422%
11	14.357	1904	1916	1922	rBV2	2369701	3719963	19.05%	2.565%
12	14.416	1922	1926	1933	rVB2	136477	223604	1.15%	0.154%
13	15.404	2087	2094	2105	rVB	131844	231189	1.18%	0.159%
14	15.851	2163	2170	2190	rBV	8536402	12551820	64.29%	8.656%
15	17.104	2377	2383	2387	rBV	3010709	4370816	22.39%	3.014%
16	17.145	2387	2390	2399	rVV	1792541	2505443	12.83%	1.728%
17	17.239	2401	2406	2410	rVB	404998	544393	2.79%	0.375%
18	17.515	2444	2453	2459	rBV3	93556	235942	1.21%	0.163%
19	17.992	2522	2534	2537	rBV2	844375	1412754	7.24%	0.974%
20	18.027	2537	2540	2546	rVV	396209	632897	3.24%	0.436%
21	18.110	2550	2554	2559	rVV	172746	265651	1.36%	0.183%
22	18.174	2559	2565	2569	rVV2	322894	634836	3.25%	0.438%
23	18.209	2569	2571	2578	rVB	171641	240224	1.23%	0.166%
24	18.486	2613	2618	2622	rBV	177542	232805	1.19%	0.161%
25	18.898	2684	2688	2695	rBV2	123763	211916	1.09%	0.146%
26	19.168	2728	2734	2742	rBV	2562886	3246212	16.63%	2.239%
27	19.274	2748	2752	2757	rBV2	317935	469621	2.41%	0.324%
28	19.498	2785	2790	2792	rBV2	442942	600160	3.07%	0.414%
29	19.533	2792	2796	2804	rVV	2087266	2852039	14.61%	1.967%
30	19.604	2804	2808	2815	rVB2	140573	218871	1.12%	0.151%
31	19.739	2823	2831	2836	rBV	16003638	19522391	100.00%	13.463%
32	19.856	2846	2851	2859	rVB3	152227	353868	1.81%	0.244%
33	20.033	2869	2881	2885	rBV3	276163	697085	3.57%	0.481%
34	20.186	2901	2907	2911	rBV2	200136	332617	1.70%	0.229%

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ClientSampleId :
TP-MH-C05

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	20.998	3039	3045	3054	rVV4	1106456	1806973	9.26%	1.246%
36	21.298	3089	3096	3100	rBV2	4007732	6047880	30.98%	4.171%
37	21.333	3100	3102	3111	rVB	903742	1028522	5.27%	0.709%
38	21.433	3116	3119	3126	rVV2	334716	464978	2.38%	0.321%
39	21.868	3190	3193	3196	rVB	355599	368348	1.89%	0.254%
40	22.327	3268	3271	3276	rBV	470889	621271	3.18%	0.428%
41	22.839	3354	3358	3360	rBV	463215	624824	3.20%	0.431%
42	22.874	3361	3364	3369	rVB	264571	525424	2.69%	0.362%
43	23.350	3442	3445	3450	rBV	212432	328939	1.68%	0.227%
44	23.409	3451	3455	3458	rVV	286225	405631	2.08%	0.280%
45	23.450	3458	3462	3471	rVB	448674	789686	4.05%	0.545%
46	23.550	3473	3479	3486	rBV2	1769448	3393178	17.38%	2.340%

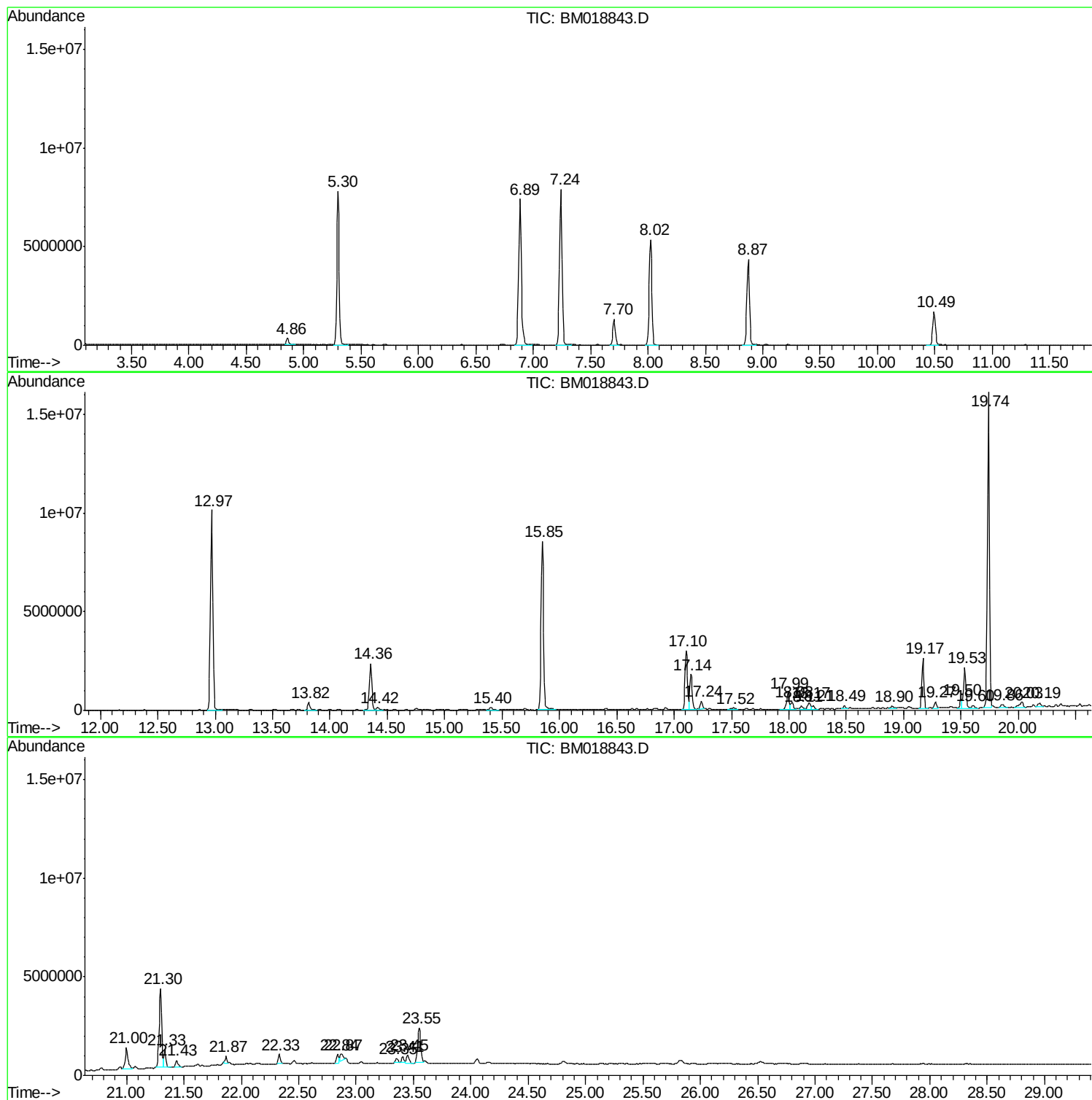
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Instrument :
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ClientSampled :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM021119.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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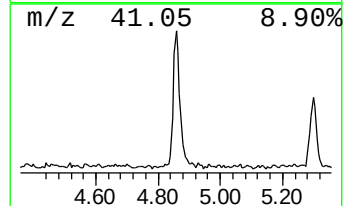
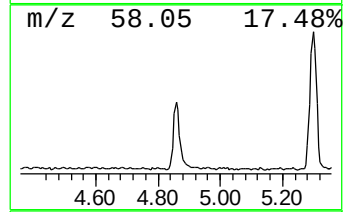
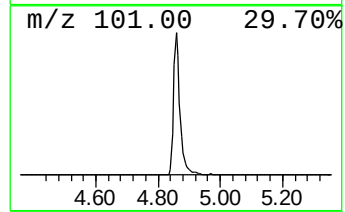
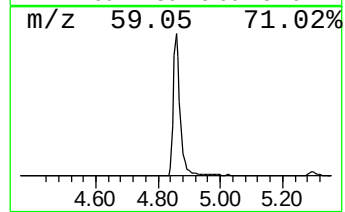
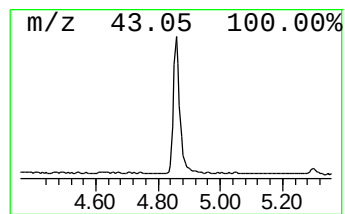
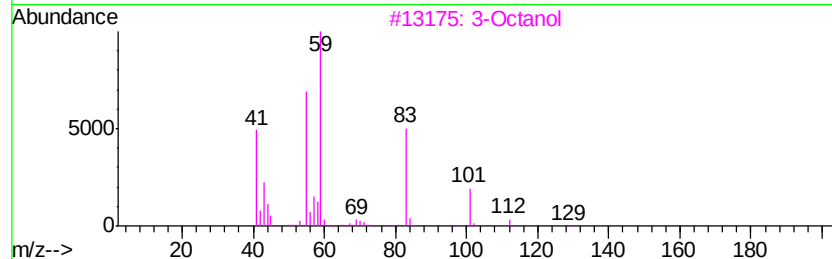
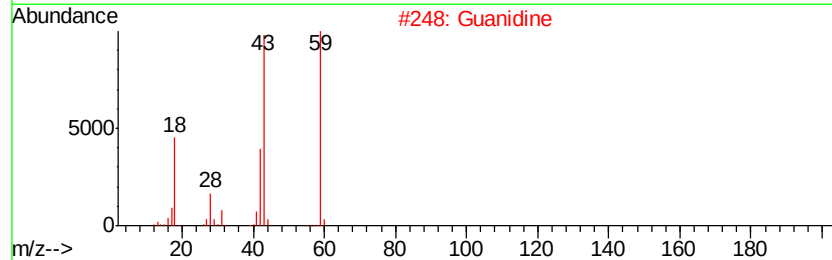
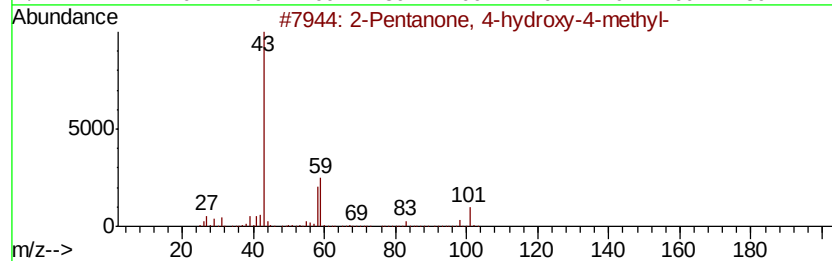
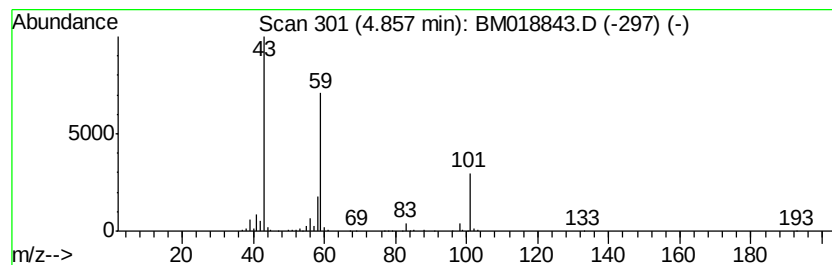
Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM021119.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
4.86	4.72 ng	486694	1,4-Dichlorobenzene-d4	7.70	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2	Guanidine	59	CH5N3	000113-00-8	37
3	3-Octanol	130	C8H18O	020296-29-1	33
4	4-Methoxy-1-pentene	100	C6H12O	098386-09-5	27
5	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25



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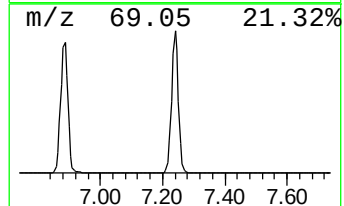
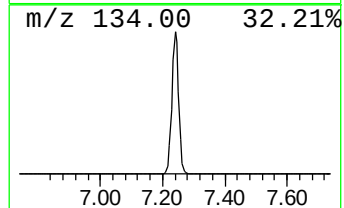
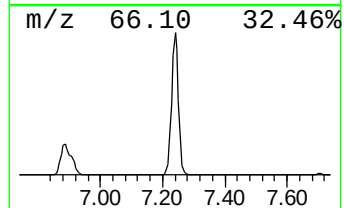
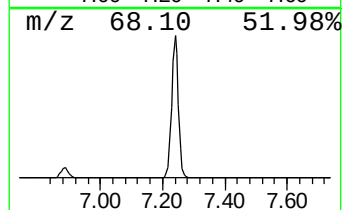
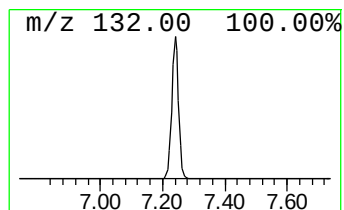
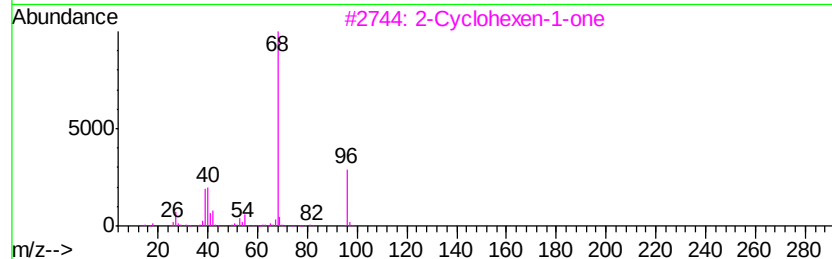
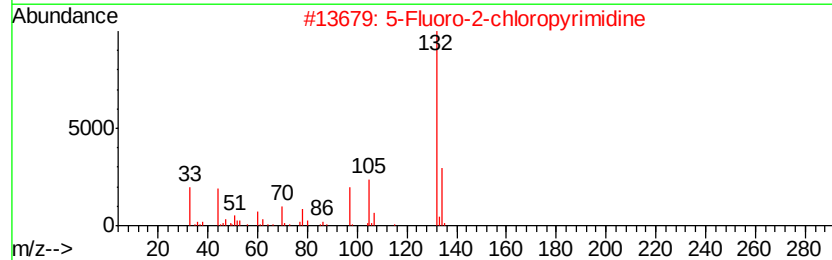
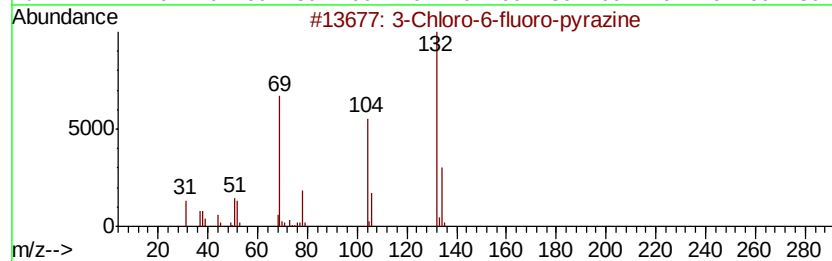
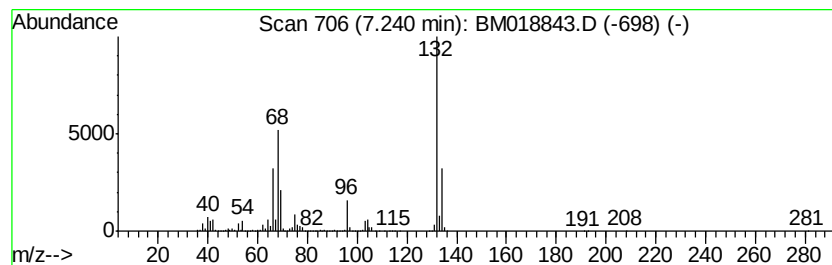
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown7.24 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.24	118.80 ng	12238300	1,4-Dichlorobenzene-d4	7.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
3		2-Cyclohexen-1-one	96	C6H8O	000930-68-7	11
4		Tranvlcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		(5-METHYL-2-PYRIDYL)ACETONITRILE	132	C8H8N2	1000241-93-9	10



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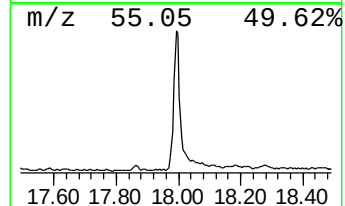
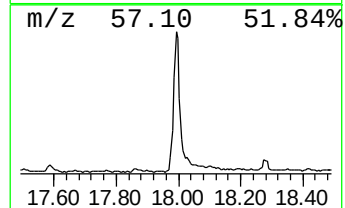
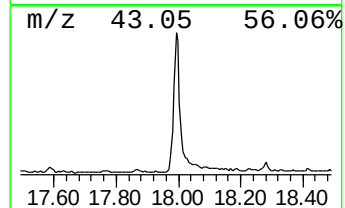
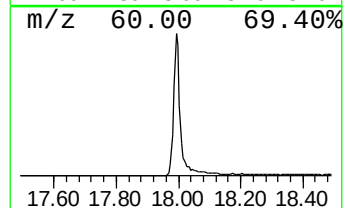
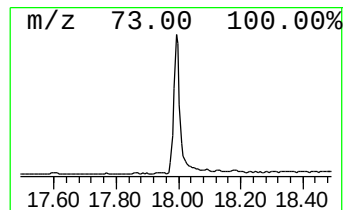
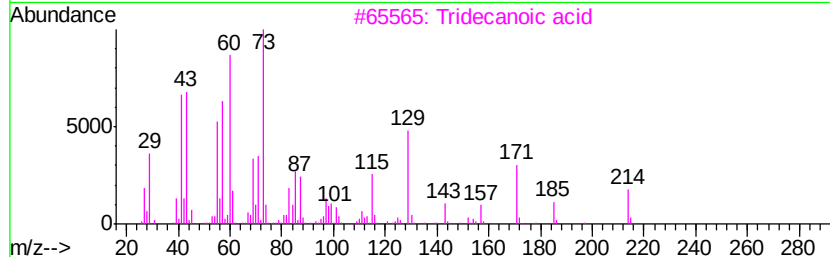
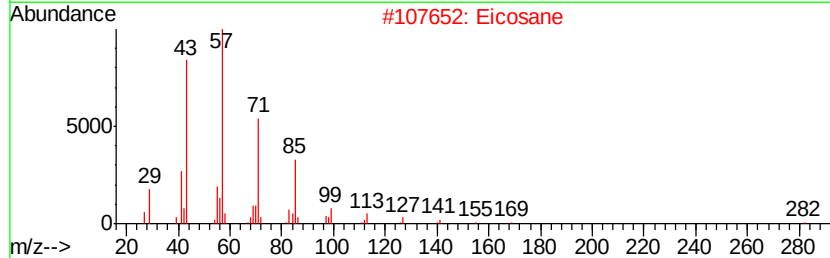
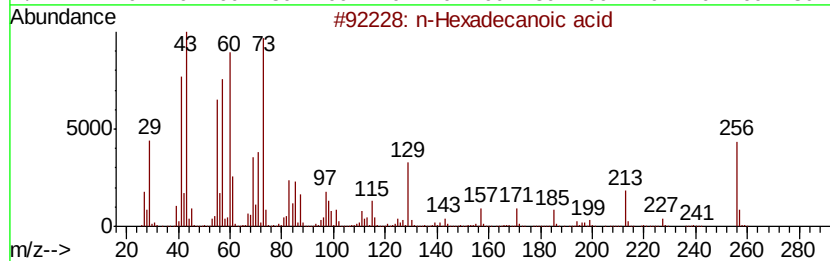
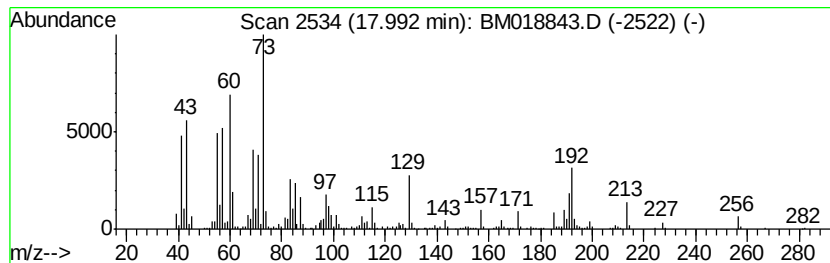
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.99	6.46 ng	1412750	Phenanthrene-d10	17.10

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2		Eicosane	282	C20H42	000112-95-8	60
3		Tridecanoic acid	214	C13H26O2	000638-53-9	50
4		n-Decanoic acid	172	C10H20O2	000334-48-5	35
5		Tetradecane	198	C14H30	000629-59-4	35



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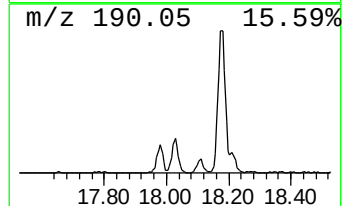
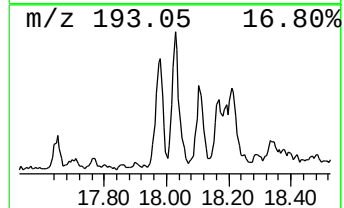
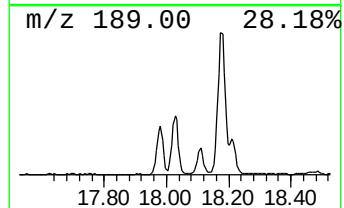
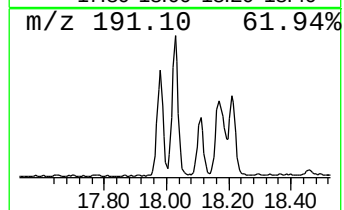
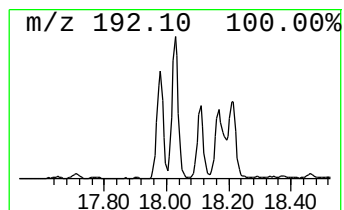
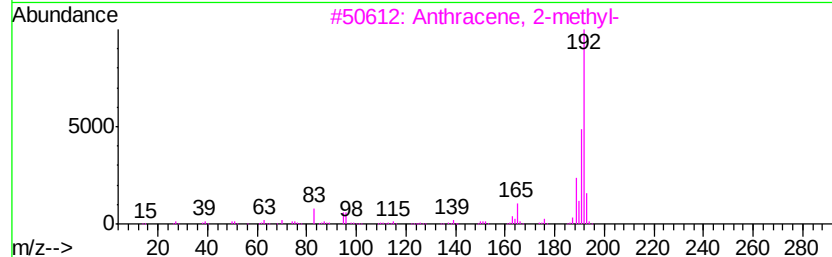
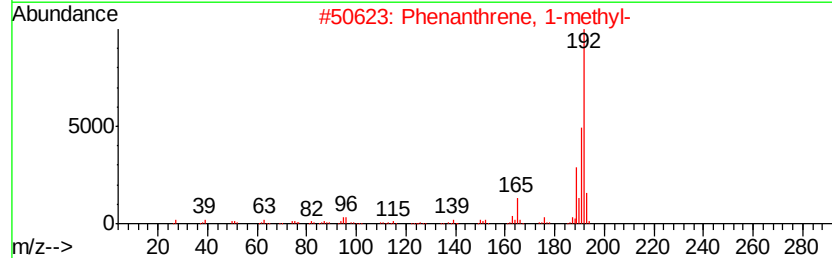
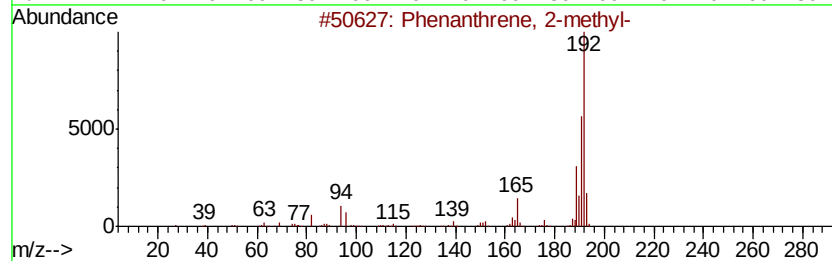
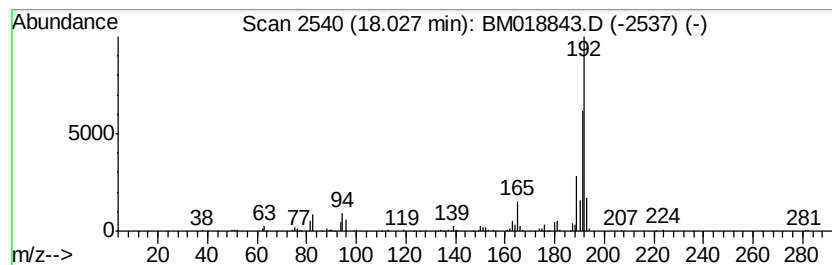
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Phenanthrene, 2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.03	2.90 ng	632897	Phenanthrene-d10	17.10	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	98
2	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	97
3	Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
4	1H-Cyclopropa[1,1]phenanthrene,1a,...	192	C15H12	000949-41-7	96
5	Anthracene, 1-methyl-	192	C15H12	000610-48-0	93



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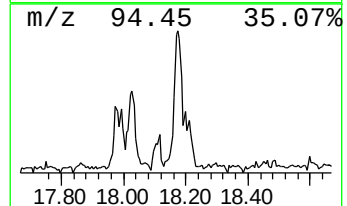
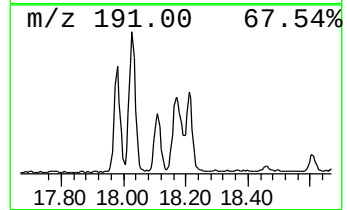
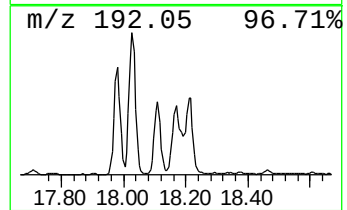
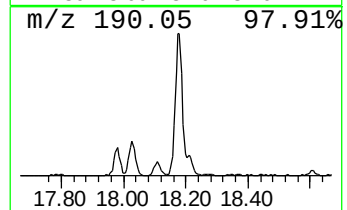
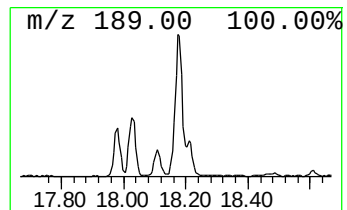
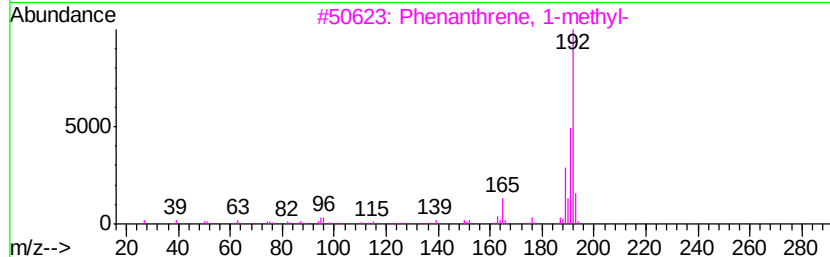
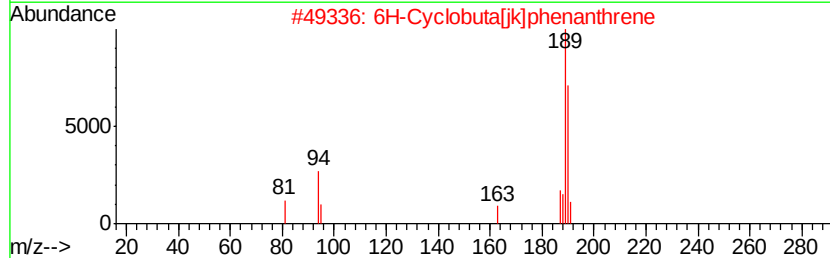
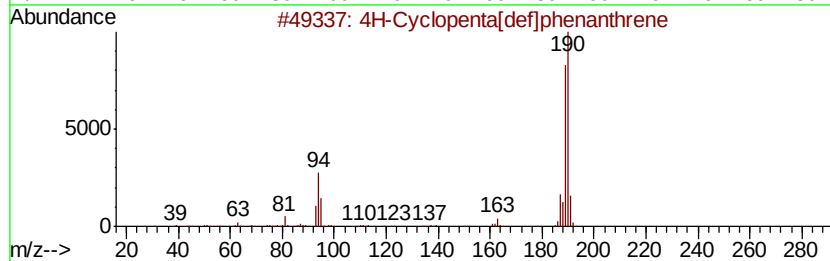
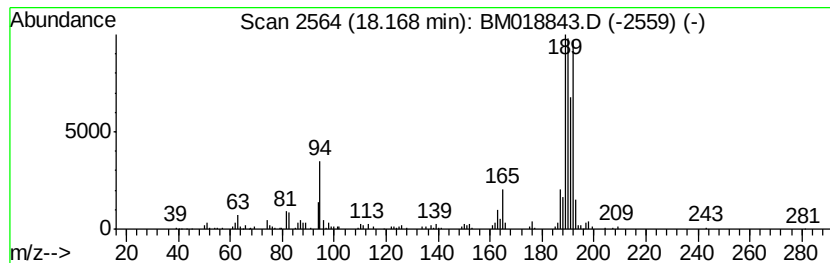
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ClientSampleId :
TP-MH-C05

Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM021119.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 4H-Cyclopenta[def]phenanthrene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.17	2.90 ng	634836	Phenanthrene-d10	17.10	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	64
2	6H-Cyclobuta[ik]phenanthrene	190	C15H10	083469-43-6	53
3	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	25
4	5H-Dibenzo[a,d]cycloheptene	192	C15H12	000256-81-5	18
5	Propyne, 1,3-diphenyl-	192	C15H12	004980-70-5	18



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM021819\
Data File : BM018843.D
Acq On : 18 Feb 2019 18:00
Operator : JU/SJ
Sample : K1518-19
Misc :
ALS Vial : 11 Sample Multiplier: 1

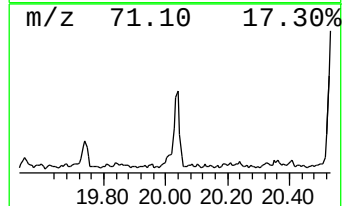
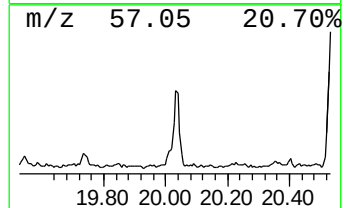
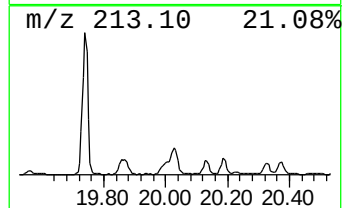
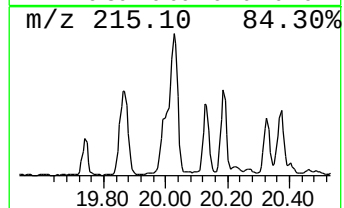
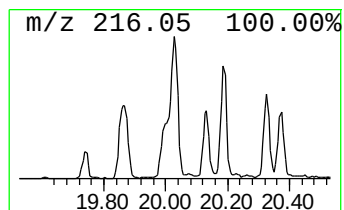
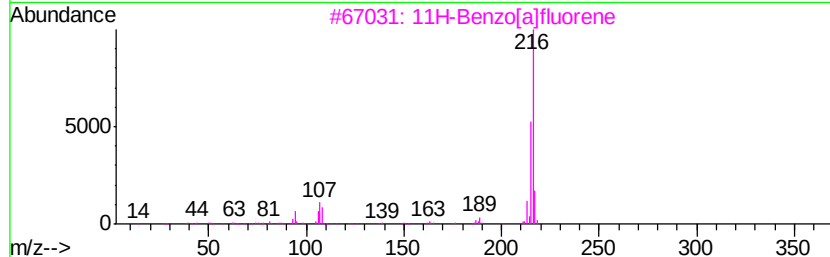
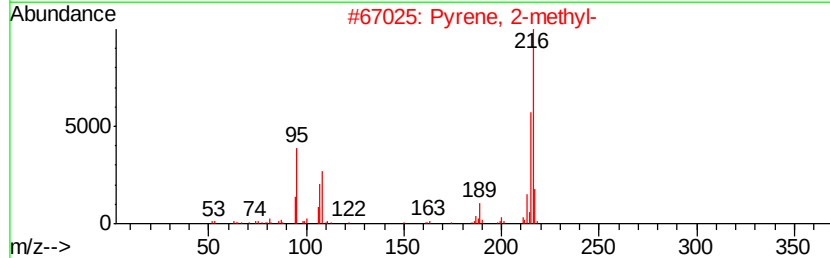
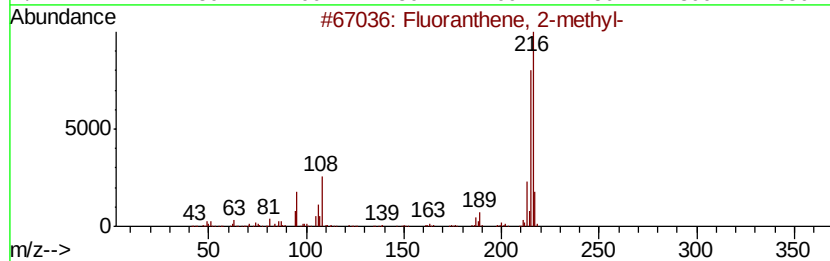
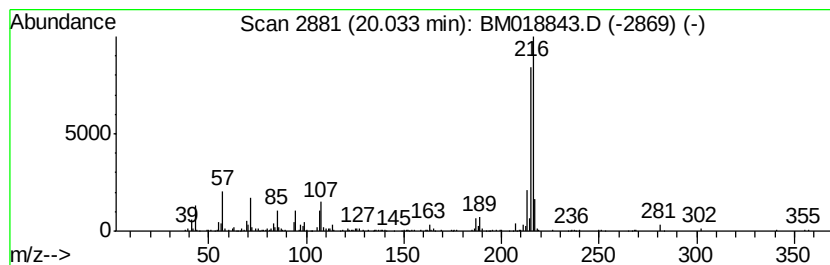
Instrument :
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ClientSampleId :
TP-MH-C05

Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM021119.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Fluoranthene, 2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.03	2.31 ng	697085	Chrysene-d12	21.30
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6 94
2	Pyrene, 2-methyl-	216	C17H12	003442-78-2 93
3	11H-Benzo[a]fluorene	216	C17H12	000238-84-6 93
4	11H-Benzo[b]fluorene	216	C17H12	000243-17-4 93
5	Pyrene, 1-methyl-	216	C17H12	002381-21-7 93



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM021819\
 Data File : BM018843.D
 Acq On : 18 Feb 2019 18:00
 Operator : JU/SJ
 Sample : K1518-19
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

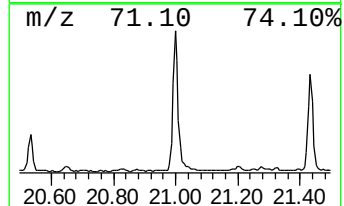
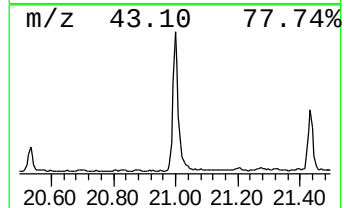
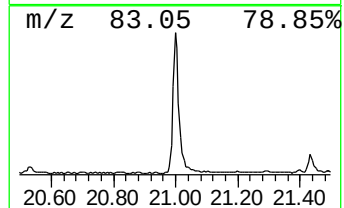
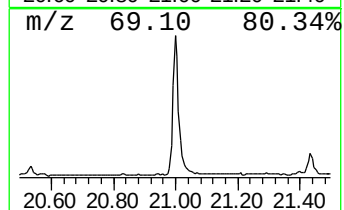
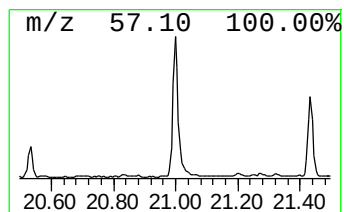
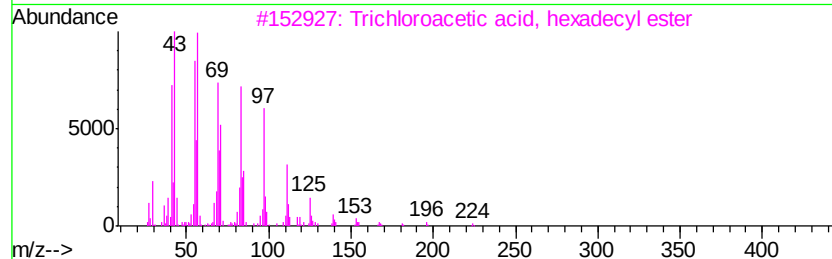
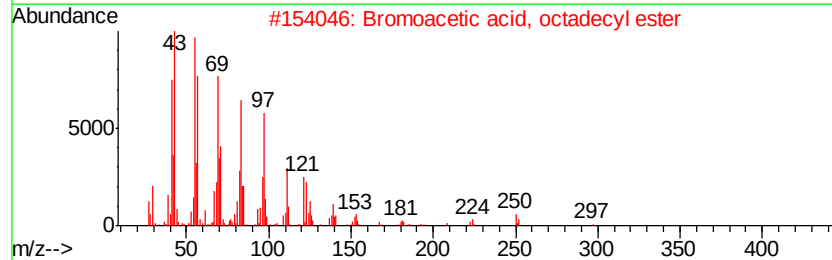
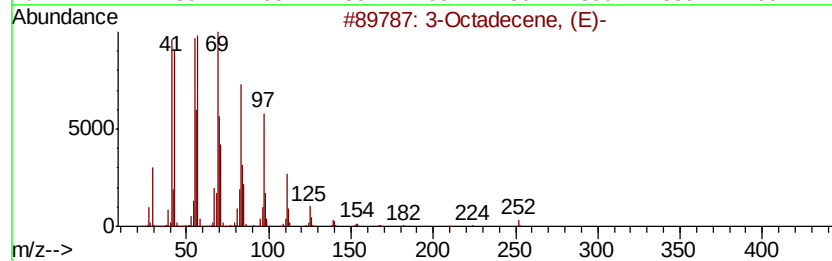
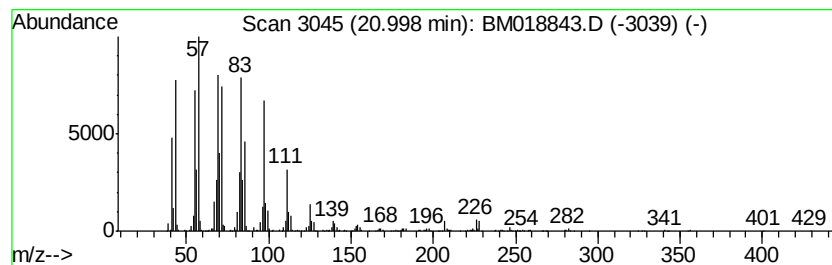
Instrument :
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 ClientSampleId :
 TP-MH-C05

Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM021119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 8 3-Octadecene, (E)- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.00	5.98 ng	1806970	Chrysene-d12	21.30
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		3-Octadecene, (E)-	252 C18H36	007206-19-1 93
2		Bromoacetic acid, octadecyl ester	390 C20H39BrO2	018992-03-5 87
3		Trichloroacetic acid, hexadecyl ...	386 C18H33Cl3O2	074339-54-1 83
4		Heptafluorobutanoic acid, heptad...	452 C21H35F7O2	1000282-97-3 83
5		Ethanol, 2-(tetradecyloxy)-	258 C16H34O2	002136-70-1 81



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM021819\
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Sample : K1518-19
Misc :
ALS Vial : 11 Sample Multiplier: 1

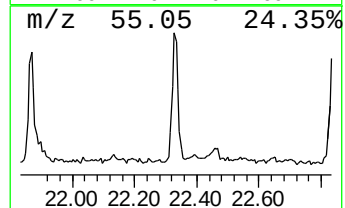
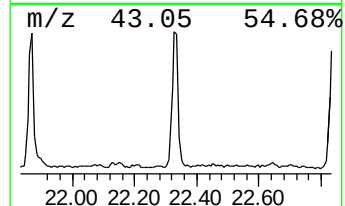
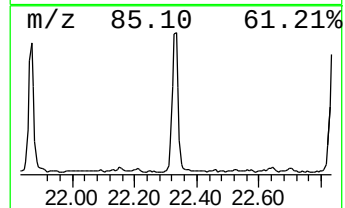
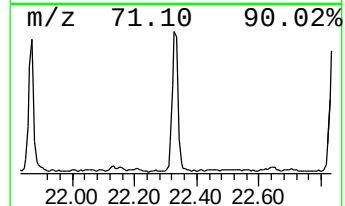
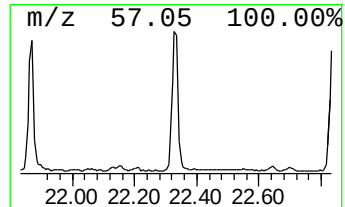
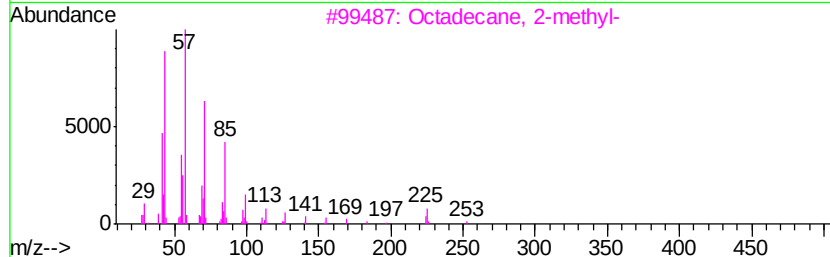
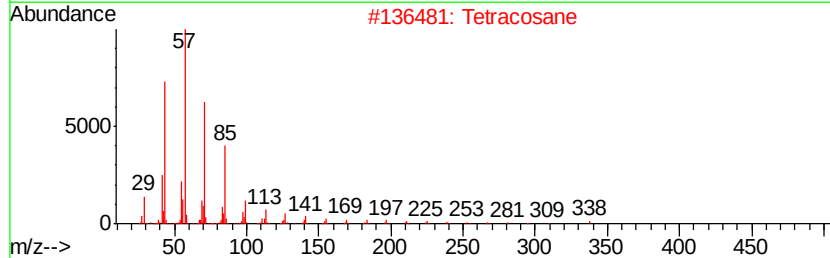
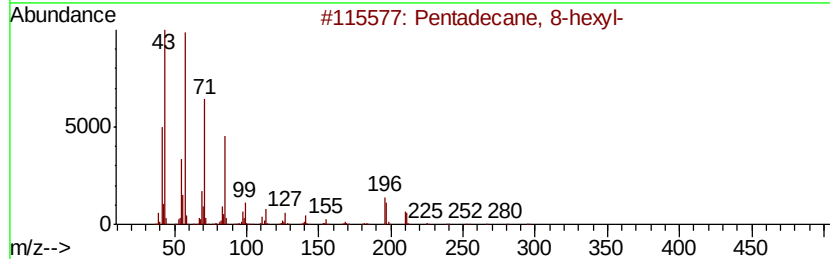
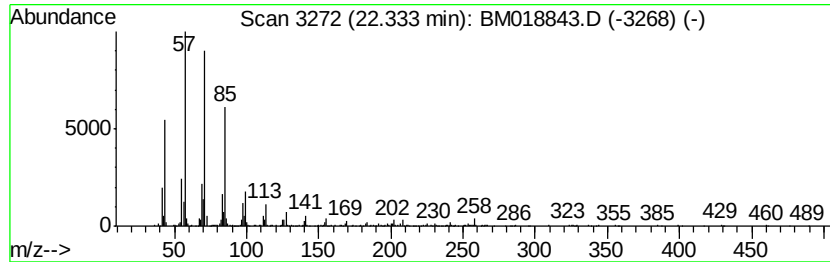
Instrument :
BNA_M
ClientSampleId :
TP-MH-C05

Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM021119.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Pentadecane, 8-hexyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.33	2.05 ng	621271	Chrysene-d12	21.30
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Pentadecane, 8-hexyl-	296 C21H44	013475-75-7	90
2	Tetracosane	338 C24H50	000646-31-1	87
3	Octadecane, 2-methyl-	268 C19H40	001560-88-9	86
4	Pentadecane, 8-heptyl-	310 C22H46	071005-15-7	86
5	Hexadecane	226 C16H34	000544-76-3	83



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Sample : K1518-19
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
TP-MH-C05

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\8270-BM021119.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...	4.86	4.7	ng	486694	1	7.70	2060320	20.0
unknown7.24	7.24	118.8	ng	12238300	1	7.70	2060320	20.0
n-Hexadecanoic acid	17.99	6.5	ng	1412750	4	17.10	4370820	20.0
Phenanthrene, 2-m...	18.03	2.9	ng	632897	4	17.10	4370820	20.0
4H-Cyclopenta[def...	18.17	2.9	ng	634836	4	17.10	4370820	20.0
Fluoranthene, 2-m...	20.03	2.3	ng	697085	5	21.30	6047880	20.0
3-Octadecene, (E)-	21.00	6.0	ng	1806970	5	21.30	6047880	20.0
Pentadecane, 8-he...	22.33	2.0	ng	621271	5	21.30	6047880	20.0