

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP120519\
 Data File : BP001224.D
 Acq On : 05 Dec 2019 21:26
 Operator : JU
 Sample : K6120-01
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 TH-7

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_P\METHODS\8270-BP112719.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	5.628	593	602	617	rBV	291543	475350	16.99%	1.748%
2	6.222	696	703	717	rBV	1845283	2300007	82.20%	8.459%
3	7.763	958	965	975	rBV	1953149	2563256	91.61%	9.427%
4	7.952	991	997	1010	rBV	2011529	2400920	85.80%	8.830%
5	8.310	1053	1058	1067	rBV	470782	555645	19.86%	2.044%
6	8.558	1094	1100	1104	rBV	1517861	1812682	64.78%	6.667%
7	9.193	1202	1208	1212	rBV	1281911	1534229	54.83%	5.643%
8	10.346	1399	1404	1409	rBV	588075	670503	23.96%	2.466%
9	11.416	1579	1586	1593	rBV2	19415	34765	1.24%	0.128%
10	12.128	1699	1707	1722	rBV	2193042	2596124	92.78%	9.548%
11	12.793	1816	1820	1825	rBV	132570	147797	5.28%	0.544%
12	13.210	1885	1891	1898	rBV	633872	814414	29.11%	2.995%
13	14.398	2086	2093	2099	rBV3	20670	40809	1.46%	0.150%
14	14.498	2104	2110	2123	rBV	1426789	1908136	68.19%	7.018%
15	14.840	2165	2168	2175	rVB2	33661	49234	1.76%	0.181%
16	15.569	2288	2292	2294	rBV	22979	30311	1.08%	0.111%
17	15.604	2294	2298	2302	rVV	728288	830231	29.67%	3.054%
18	15.640	2302	2304	2310	rVB	42737	46969	1.68%	0.173%
19	15.745	2319	2322	2327	rVB6	28021	32017	1.14%	0.118%
20	16.492	2446	2449	2455	rBV2	130483	176182	6.30%	0.648%
21	17.351	2591	2595	2601	rVB	71512	87358	3.12%	0.321%
22	17.581	2631	2634	2640	rBV3	37759	55675	1.99%	0.205%
23	17.651	2643	2646	2651	rVB	76955	90308	3.23%	0.332%
24	17.786	2667	2669	2673	rVB	38100	37827	1.35%	0.139%
25	17.892	2681	2687	2691	rBV	2449973	2798125	100.00%	10.291%
26	18.263	2746	2750	2753	rVB	74446	85069	3.04%	0.313%
27	18.710	2822	2826	2830	rBV	119437	129803	4.64%	0.477%
28	19.086	2887	2890	2893	rBV	58391	69094	2.47%	0.254%
29	19.128	2893	2897	2912	rBV2	308075	626287	22.38%	2.303%
30	19.245	2912	2917	2921	rBV	703962	818020	29.23%	3.009%
31	19.539	2964	2967	2970	rBV	159427	162417	5.80%	0.597%
32	19.933	3030	3034	3036	rBV	214607	278308	9.95%	1.024%
33	20.116	3062	3065	3069	rVB	221922	240306	8.59%	0.884%
34	20.304	3095	3097	3101	rVB	172050	161674	5.78%	0.595%

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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

Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP112719.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	20.704	3161	3165	3170	rVB	237069	318980	11.40%	1.173%
36	21.022	3215	3219	3227	rVB	586883	916357	32.75%	3.370%
37	21.145	3236	3240	3246	rVB	186797	275484	9.85%	1.013%
38	21.645	3320	3325	3344	rVB4	274594	812656	29.04%	2.989%
39	22.216	3419	3422	3429	rVB2	123186	206065	7.36%	0.758%

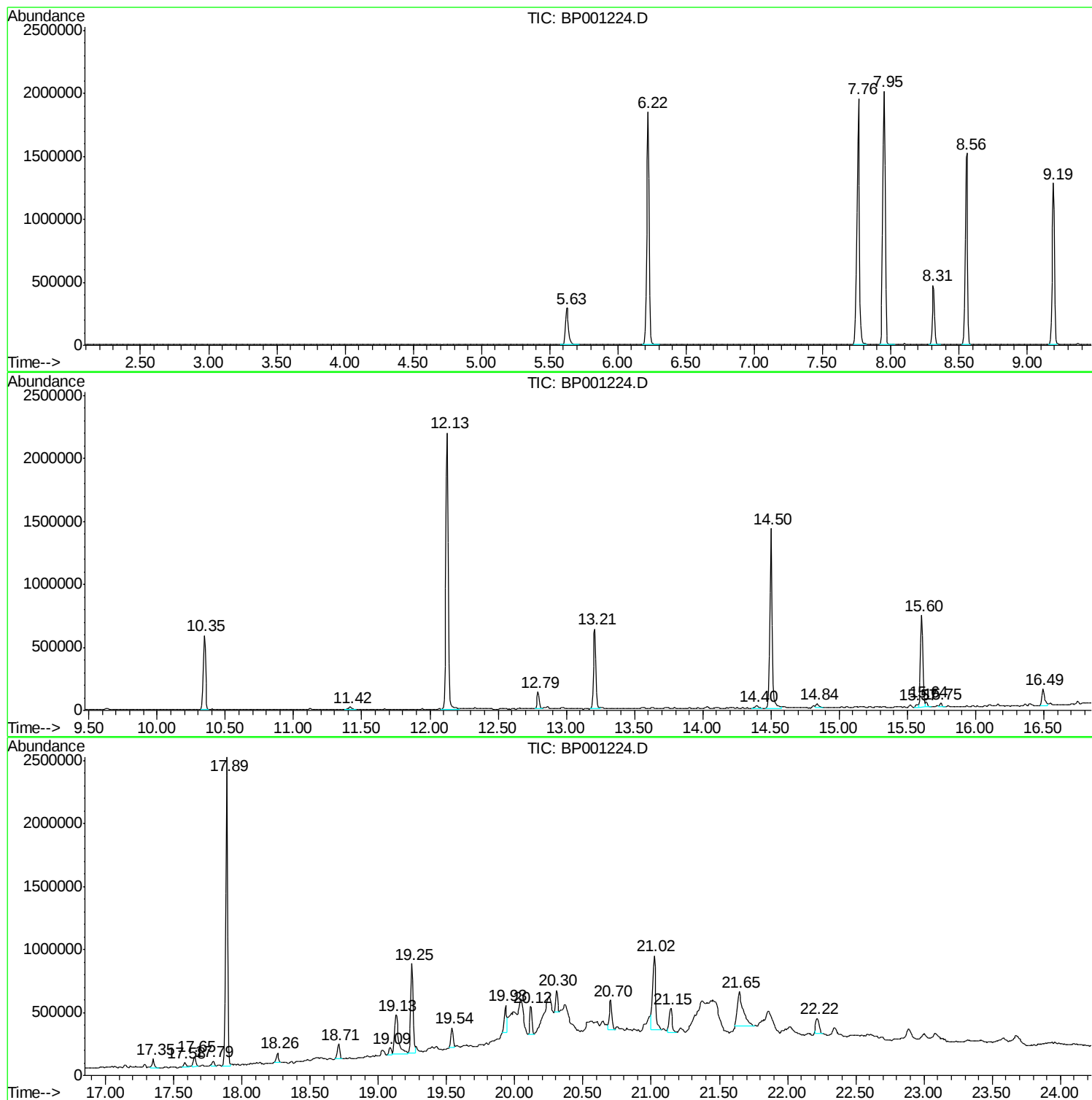
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP112719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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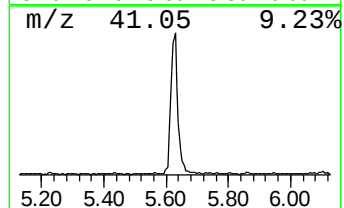
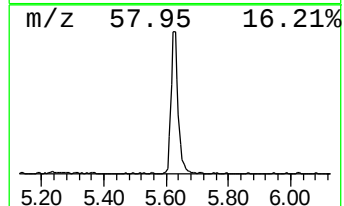
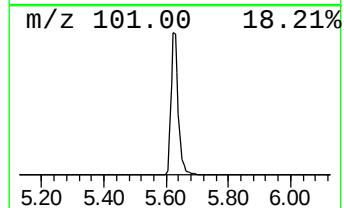
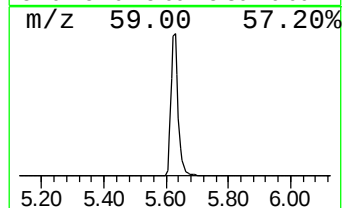
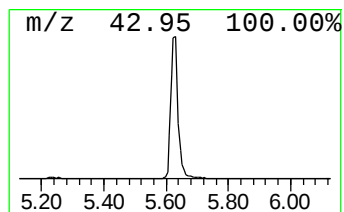
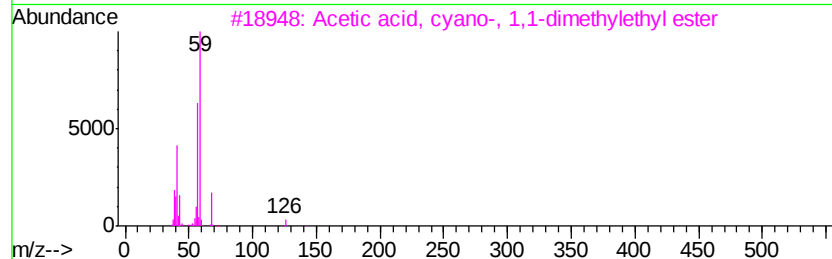
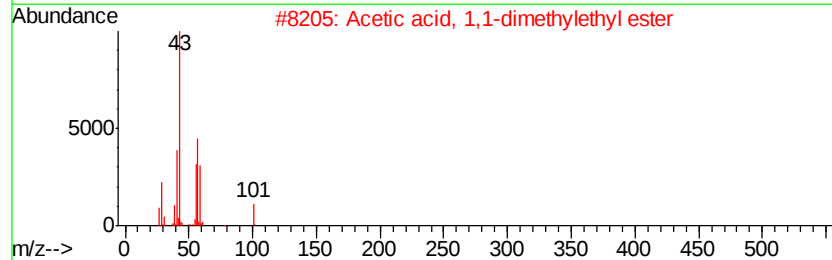
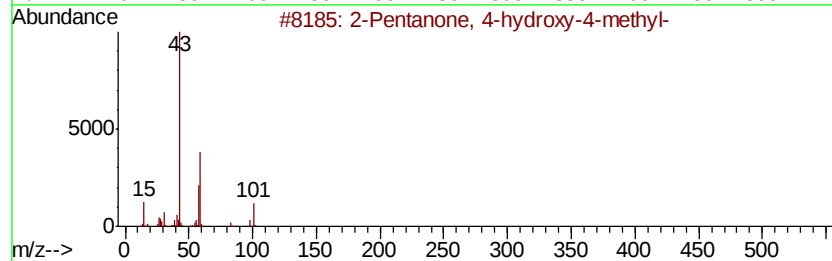
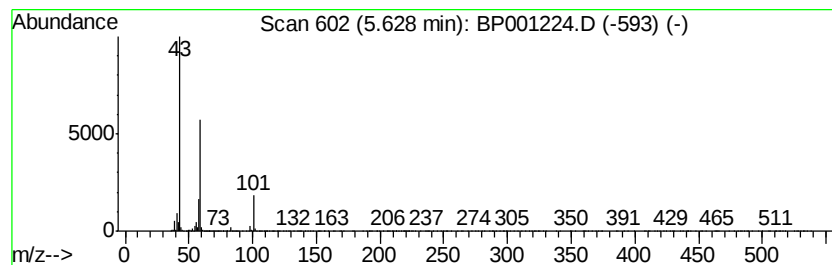
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.63	17.11 ng	475350	1,4-Dichlorobenzene-d4	8.31

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-			116	C6H12O2	000123-42-2	59
2	Acetic acid, 1,1-dimethylethyl e...			116	C6H12O2	000540-88-5	28
3	Acetic acid, cyano-, 1,1-dimethyl...			141	C7H11NO2	001116-98-9	23
4	1-Propen-2-ol, acetate			100	C5H8O2	000108-22-5	10
5	Morpholine, 4-methyl-			101	C5H11NO	000109-02-4	9



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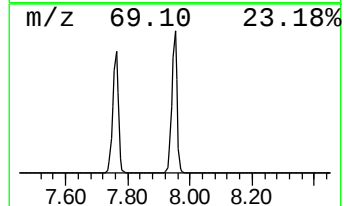
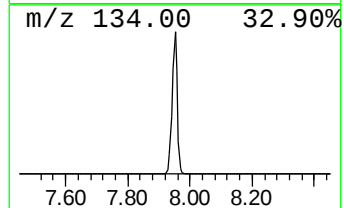
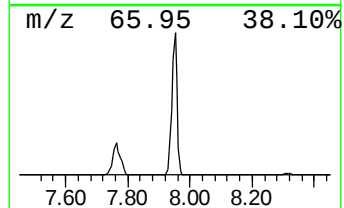
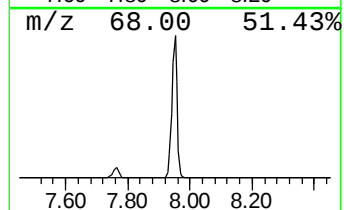
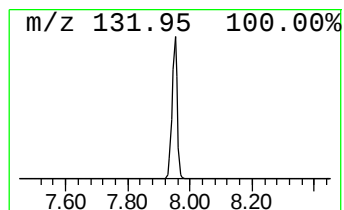
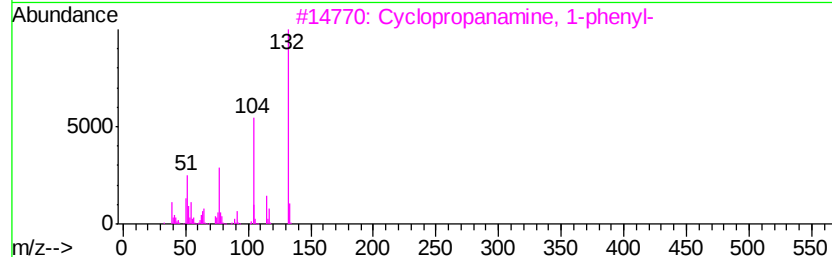
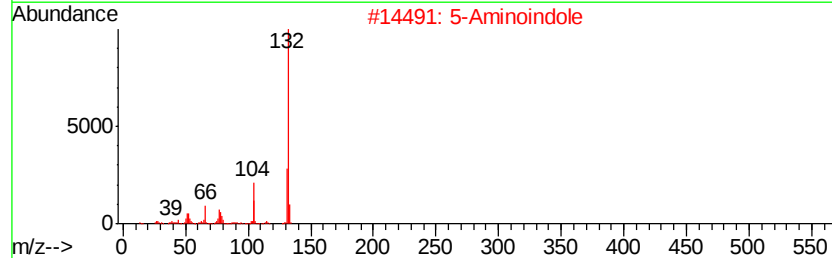
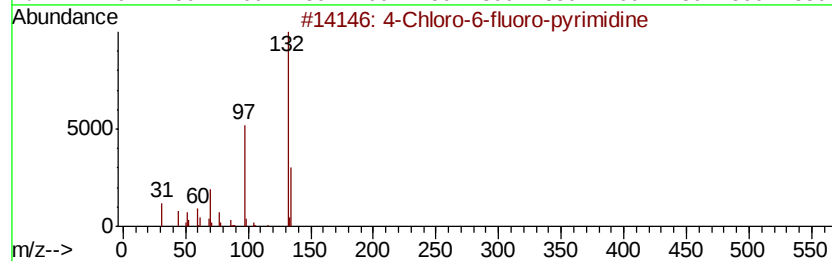
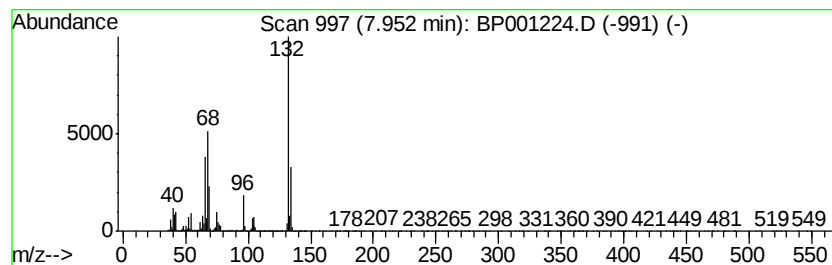
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 2 unknown7.95 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.95	86.42 ng	2400920	1,4-Dichlorobenzene-d4	8.31

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	14
2		5-Aminoindole	132	C8H8N2	005192-03-0	10
3		Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	10
4		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	10
5		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	10



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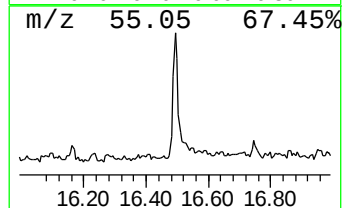
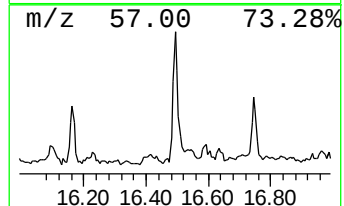
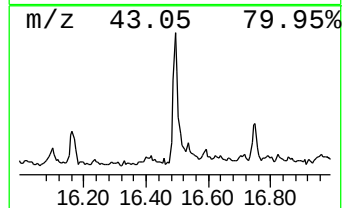
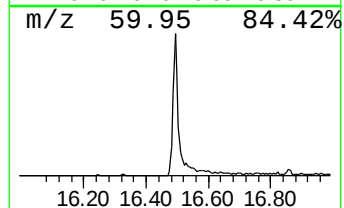
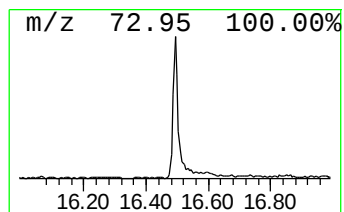
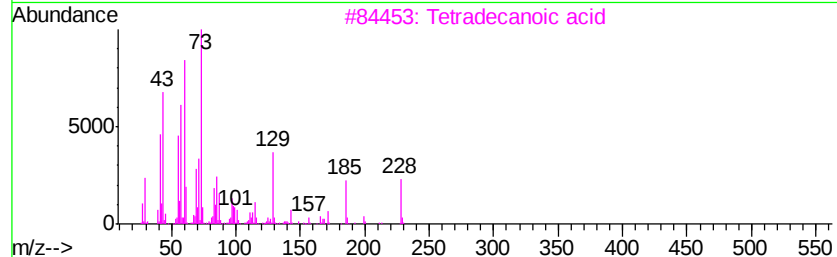
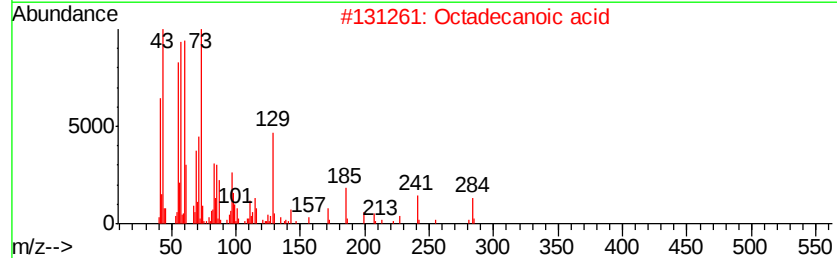
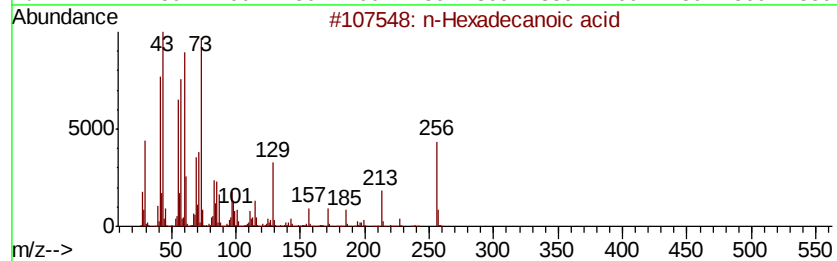
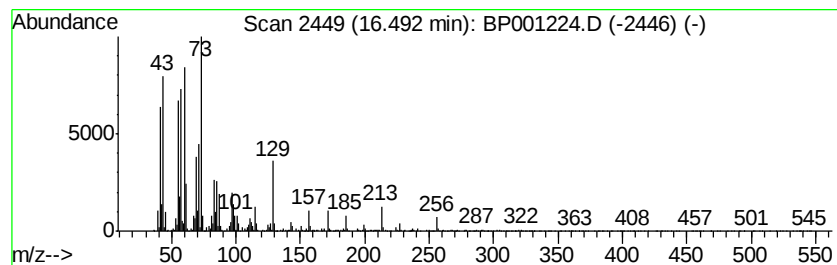
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 n-Hexadecanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD		R.T.
16.49	4.24 ng	176182	Phenanthrene-d10		15.60
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2	Octadecanoic acid	284	C18H36O2	000057-11-4	90
3	Tetradecanoic acid	228	C14H28O2	000544-63-8	86
4	Pentadecanoic acid	242	C15H30O2	001002-84-2	83
5	Tridecanoic acid	214	C13H26O2	000638-53-9	68



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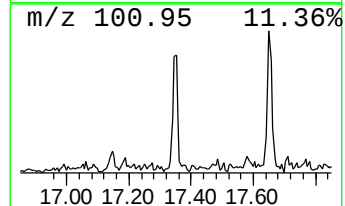
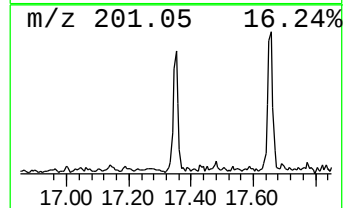
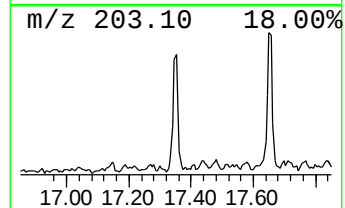
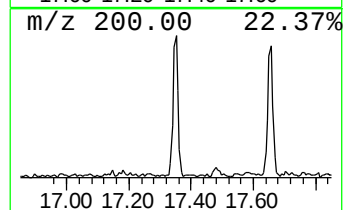
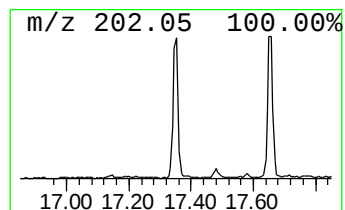
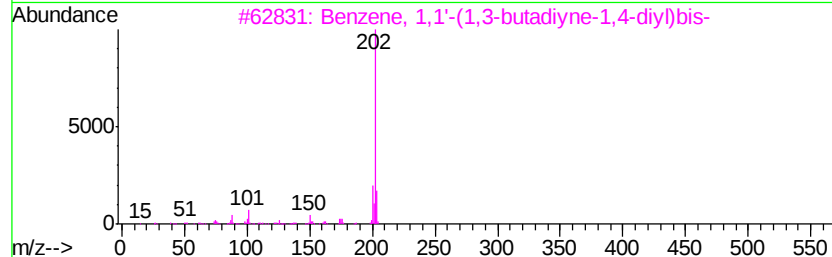
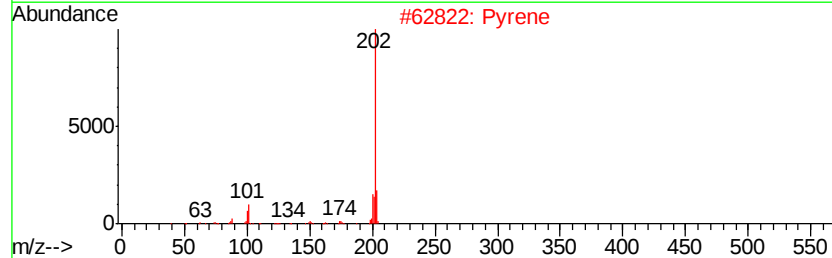
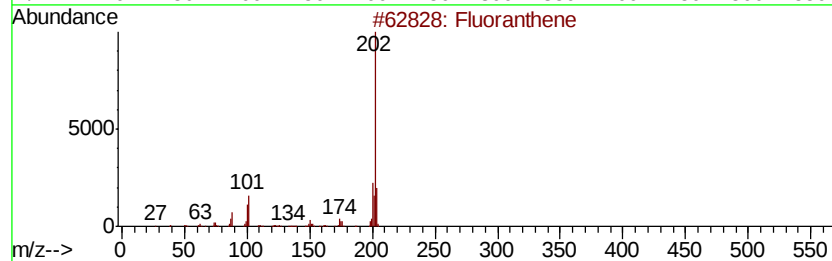
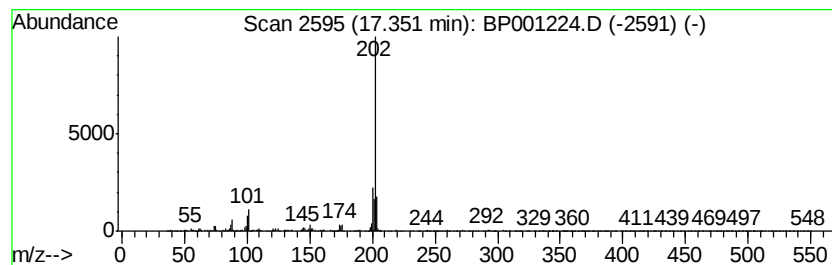
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 5 Benzene, 1,1'-(1,3-butadiyn... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.35	2.10 ng	87358	Phenanthrene-d10	15.60

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Fluoranthene	202	C16H10	000206-44-0	98
2		Pyrene	202	C16H10	000129-00-0	81
3		Benzene, 1,1'-(1,3-butadiyne-1,4...	202	C16H10	000886-66-8	81
4		1,4-Pentadiyn-3-one, 1,5-diphenyl-	230	C17H10O	015814-30-9	64
5		3H-Benzo[4,5]furo[3,2-d]pyrimidi...	202	C10H6N2OS	062208-70-2	50



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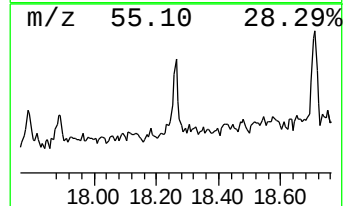
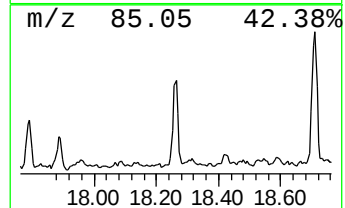
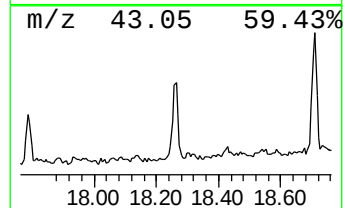
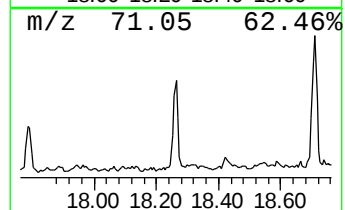
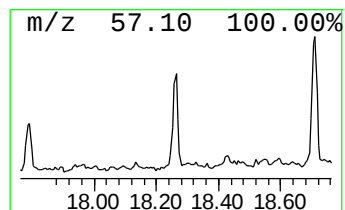
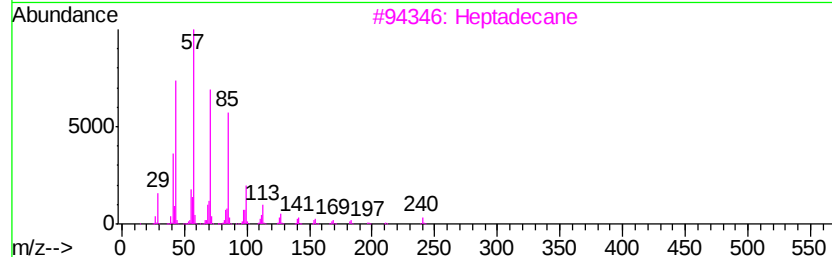
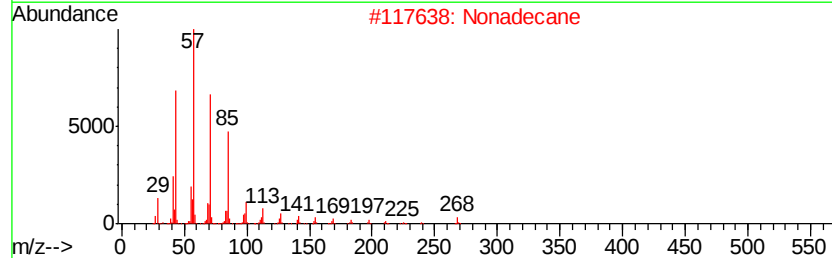
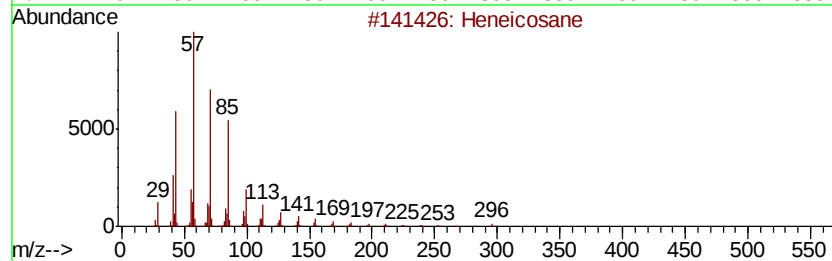
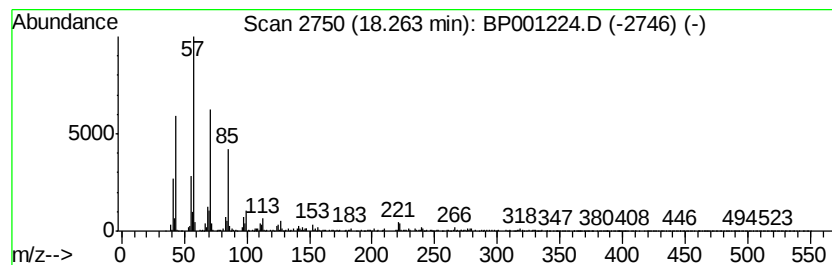
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TIC Integration Parameters: LSCINT.P

Peak Number 7 Heneicosane Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.26	2.08 ng	85069	Chrysene-d12	19.25

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	95
2		Nonadecane	268	C19H40	000629-92-5	95
3		Heptadecane	240	C17H36	000629-78-7	94
4		Docosane	310	C22H46	000629-97-0	93
5		Heptacosane	380	C27H56	000593-49-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP120519\
Data File : BP001224.D
Acq On : 05 Dec 2019 21:26
Operator : JU
Sample : K6120-01
Misc :
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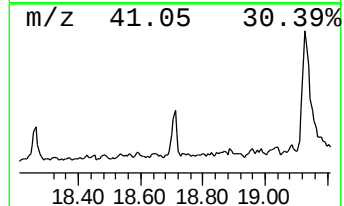
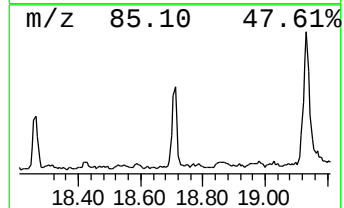
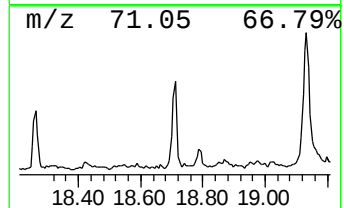
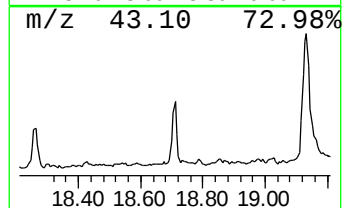
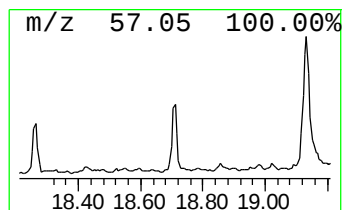
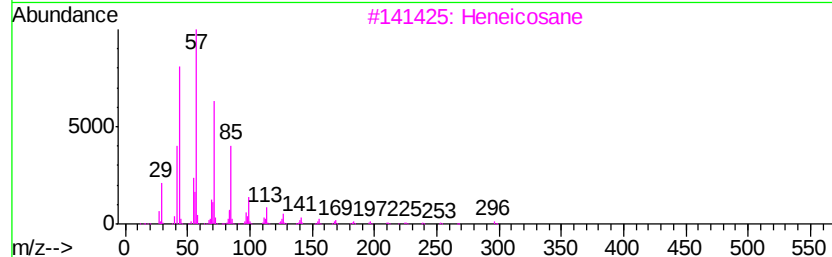
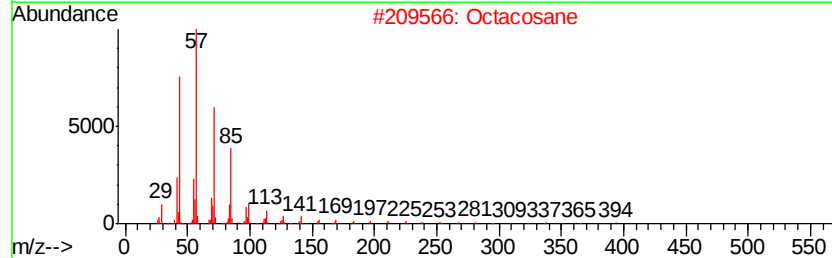
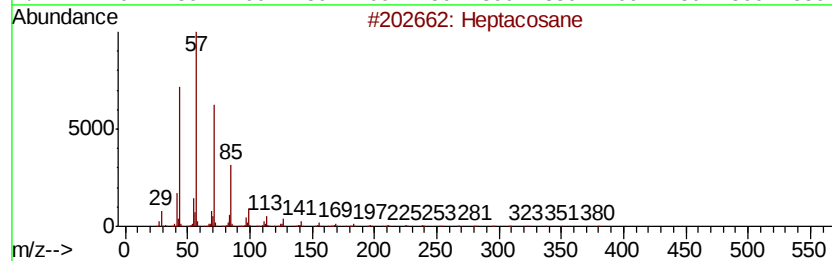
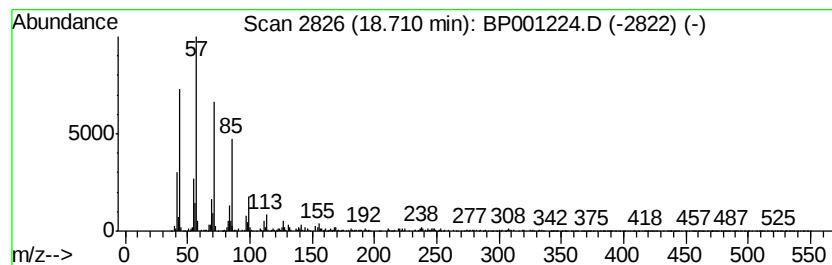
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP112719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Heptacosane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD		R.T.
18.71	3.17 ng	129803	Chrysene-d12		19.25
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptacosane	380	C27H56	000593-49-7	91
2	Octacosane	394	C28H58	000630-02-4	87
3	Heneicosane	296	C21H44	000629-94-7	83
4	1-Iodoundecane	282	C11H23I	004282-44-4	83
5	Hexatriacontane	507	C36H74	000630-06-8	80



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP120519\
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Instrument :
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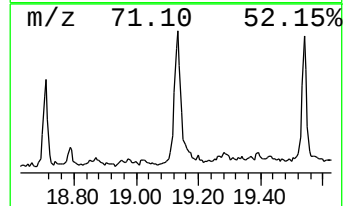
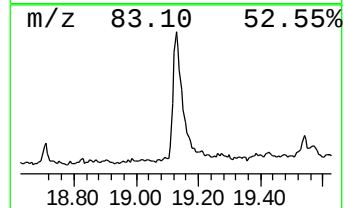
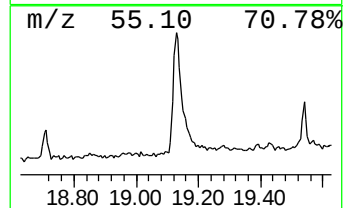
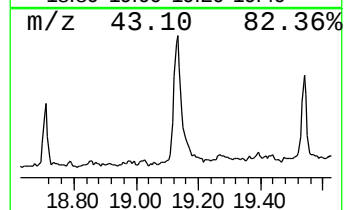
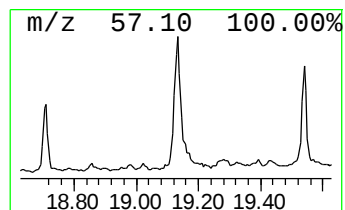
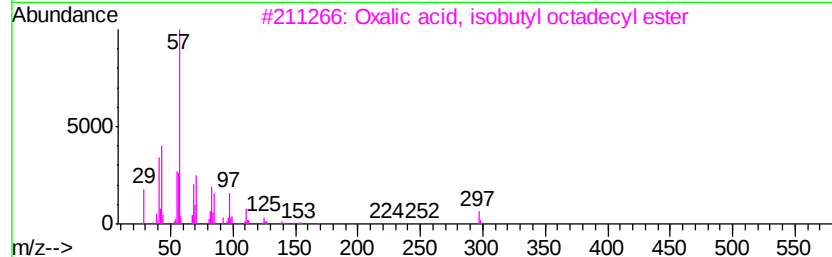
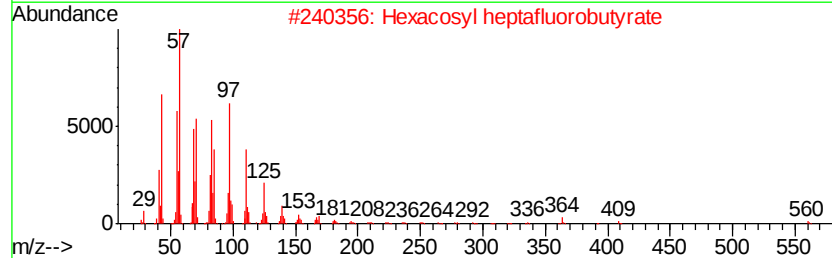
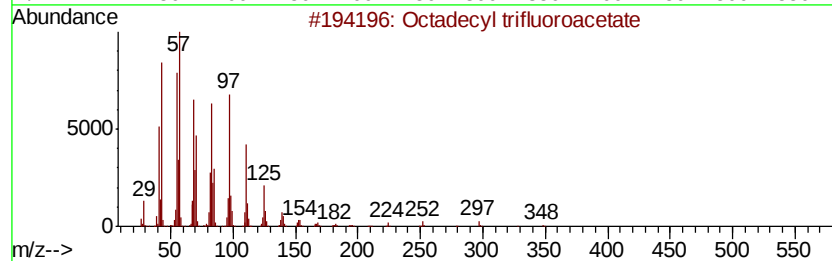
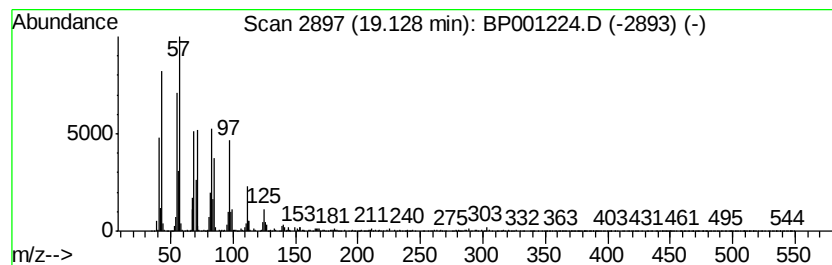
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Octadecyl trifluoroacetate Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.13	15.31 ng	626287	Chrysene-d12	19.25

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octadecyl trifluoroacetate	366	C20H37F3O2	079392-43-1	91
2		Hexacosyl heptafluorobutyrate	578	C30H53F7O2	1000351-83-3	91
3		Oxalic acid, isobutyl octadecyl ...	398	C24H46O4	1000309-38-3	91
4		Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	91
5		Tricosyl trifluoroacetate	436	C25H47F3O2	1000351-75-1	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP120519\
Data File : BP001224.D
Acq On : 05 Dec 2019 21:26
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Misc :
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Instrument :
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ClientSampleId :
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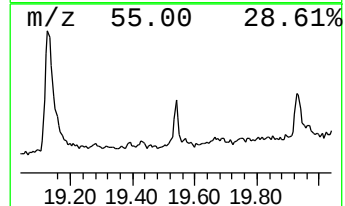
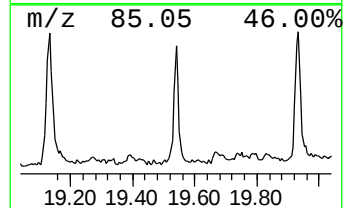
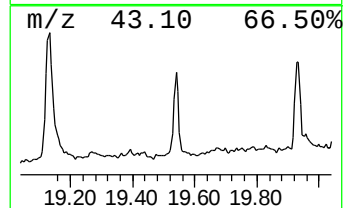
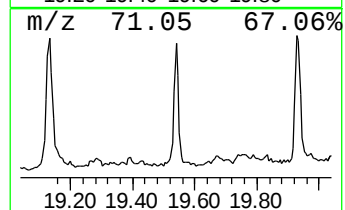
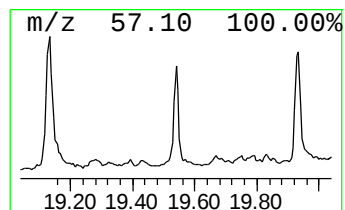
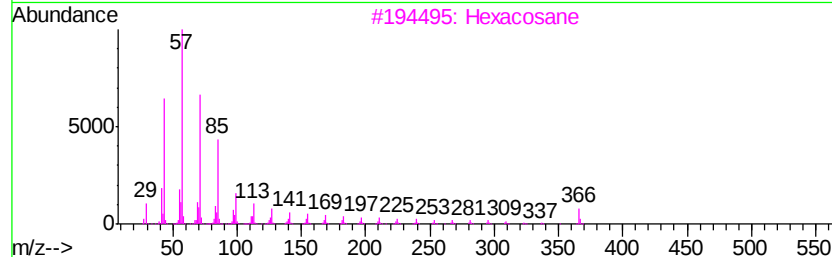
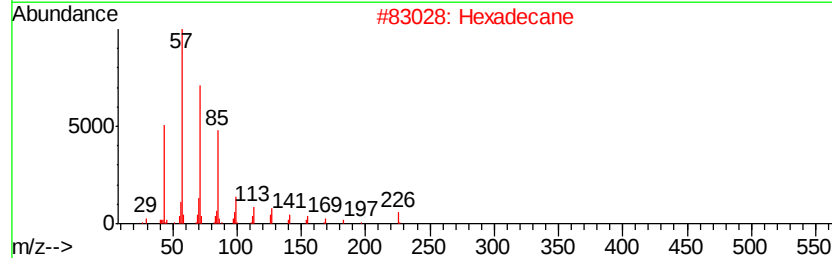
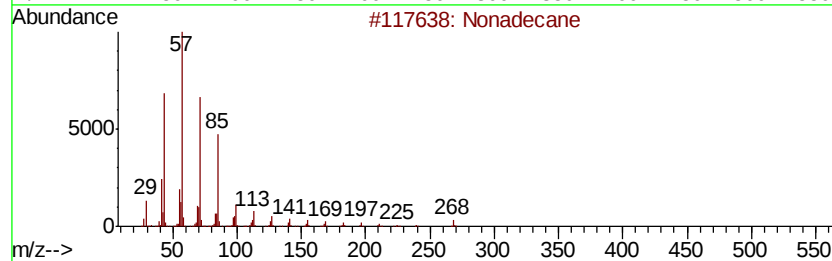
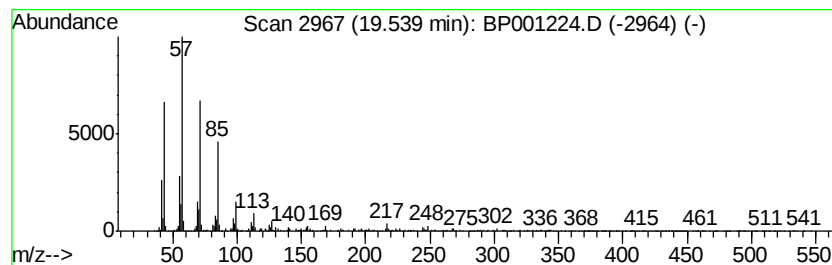
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Nonadecane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.54	3.97 ng	162417	Chrysene-d12	19.25

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonadecane	268	C19H40	000629-92-5	93
2		Hexadecane	226	C16H34	000544-76-3	93
3		Hexacosane	366	C26H54	000630-01-3	90
4		Heneicosane	296	C21H44	000629-94-7	90
5		Docosane, 11-butyl-	366	C26H54	013475-76-8	89



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP120519\
Data File : BP001224.D
Acq On : 05 Dec 2019 21:26
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Misc :
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Instrument :
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ClientSampleId :
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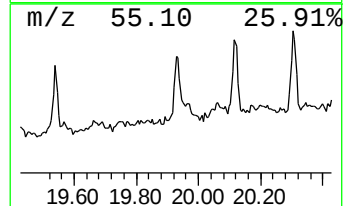
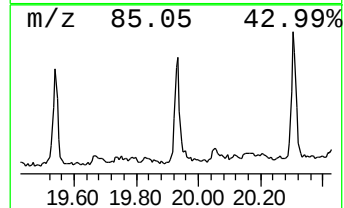
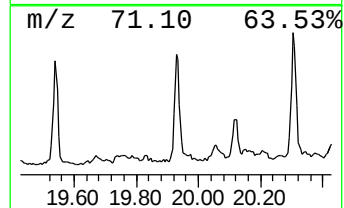
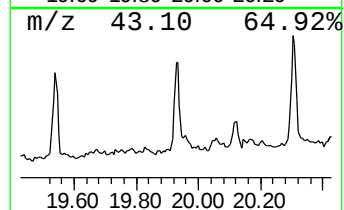
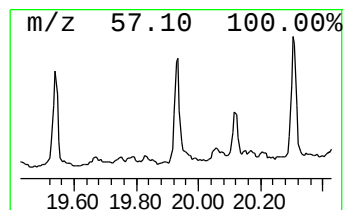
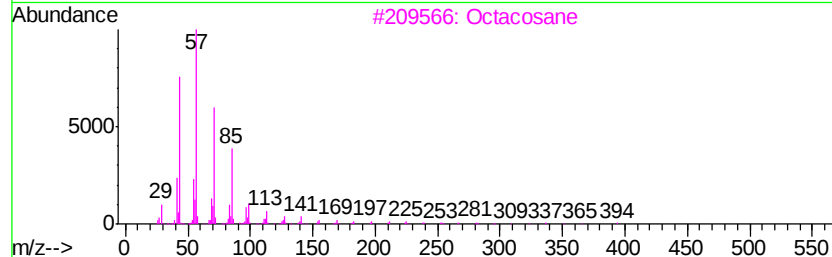
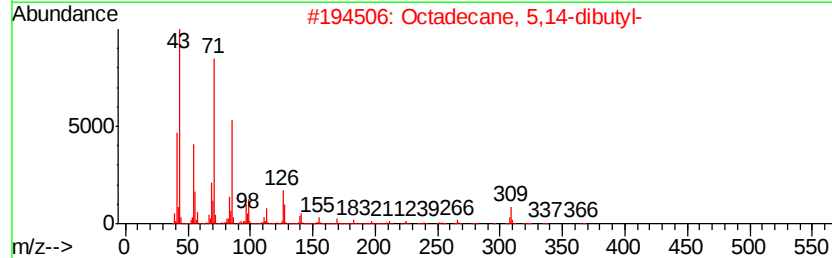
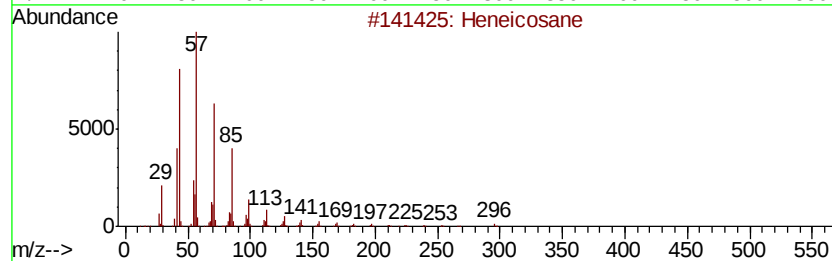
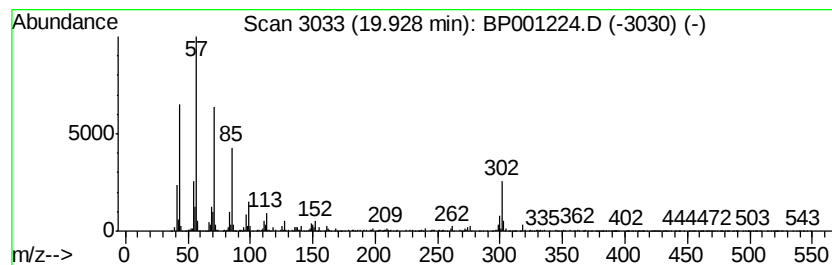
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Octadecane, 5,14-dibutyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.93	6.80 ng	278308	Chrysene-d12	19.25

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	95
2			Octadecane, 5,14-dibutyl-	366	C26H54	055282-13-8	91
3			Octacosane	394	C28H58	000630-02-4	70
4			Tridecane, 1-iodo-	310	C13H27I	035599-77-0	64
5			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	62



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP120519\
 Data File : BP001224.D
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Instrument :
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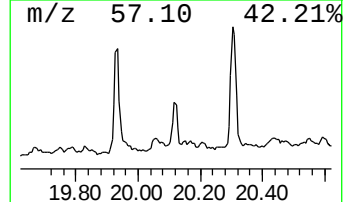
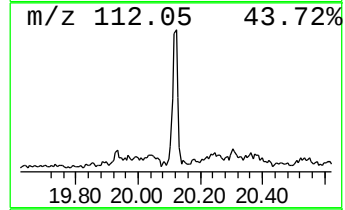
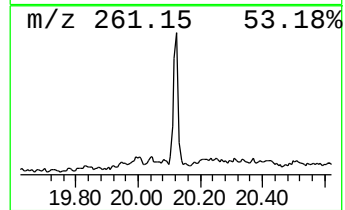
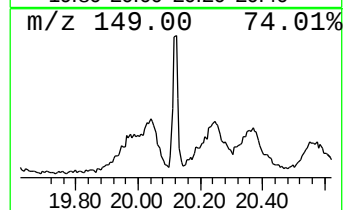
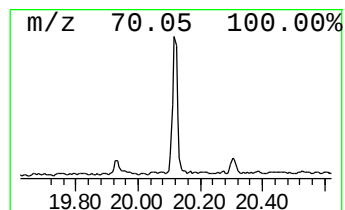
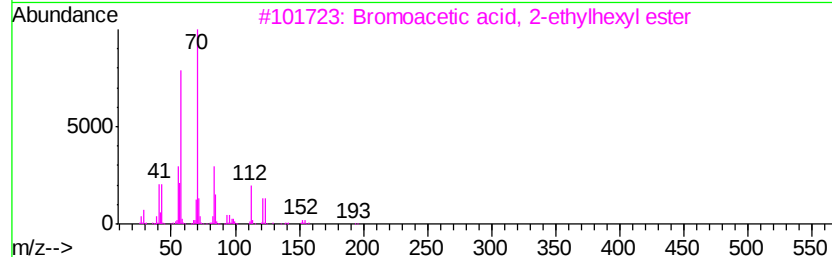
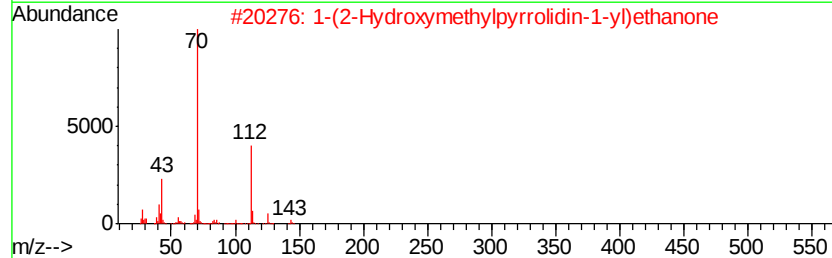
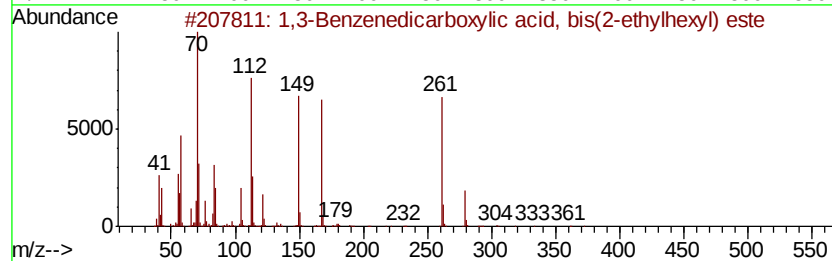
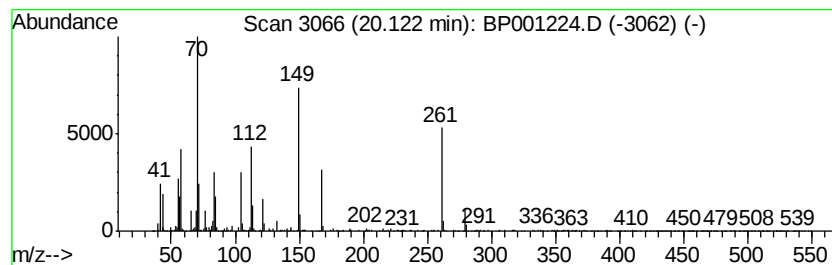
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 12 unknown20.12 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.12	5.88 ng	240306	Chrysene-d12	19.25

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,3-Benzenedicarboxylic acid, bi...	390	C24H38O4	000137-89-3	38
2			1-(2-Hydroxymethylpyrrolidin-1-yl)...	143	C7H13NO2	1000192-83-2	30
3			Bromoacetic acid, 2-ethylhexyl e...	250	C10H19BrO2	068144-73-0	25
4			2-Propenoic acid, 2-methyl-, oct...	198	C12H22O2	002157-01-9	23
5			Fumaric acid, 2-ethylhexyl hexyl...	312	C18H32O4	1000339-14-1	22



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP120519\
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Acq On : 05 Dec 2019 21:26
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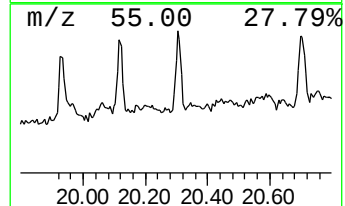
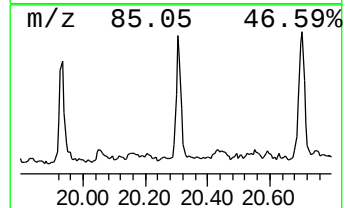
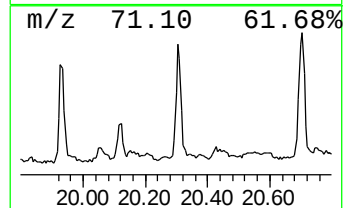
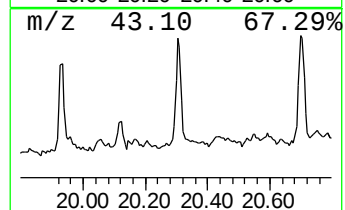
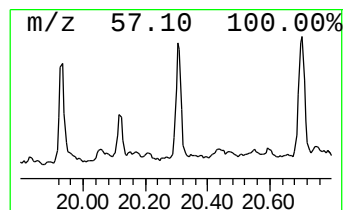
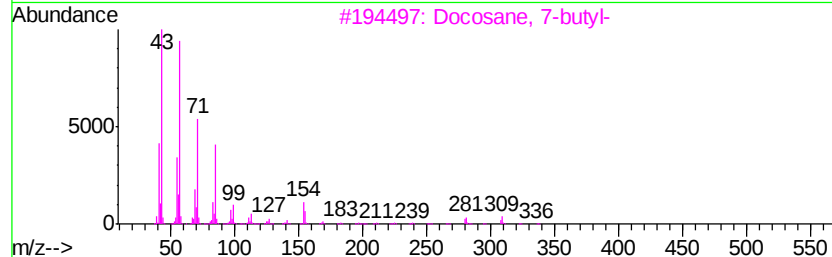
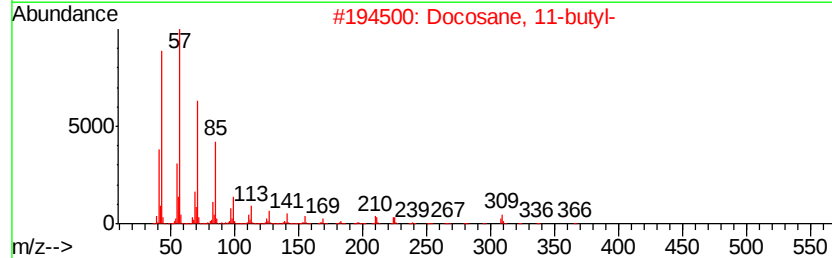
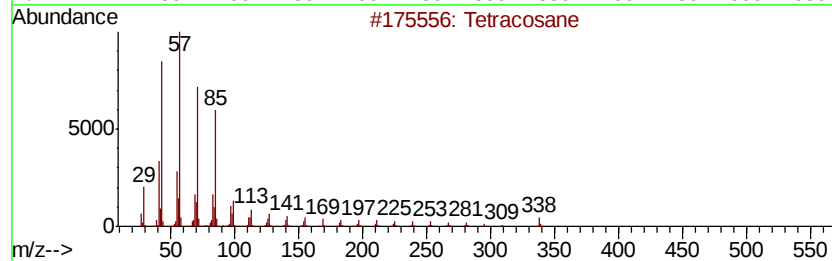
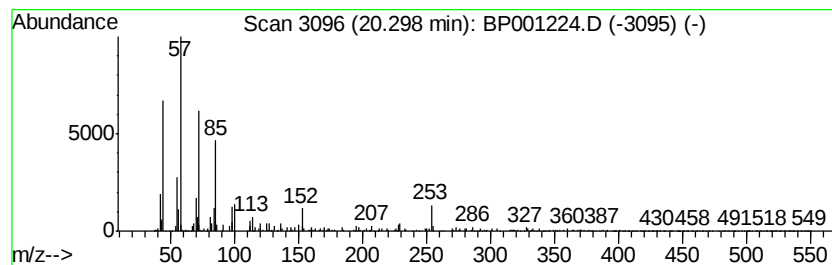
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Tetracosane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.	
20.30	3.53 ng	161674	Perylene-d12	21.02	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetracosane	338	C24H50	000646-31-1	91
2	Docosane, 11-butyl-	366	C26H54	013475-76-8	80
3	Docosane, 7-butyl-	366	C26H54	055282-15-0	80
4	Tridecane, 7-propyl-	226	C16H34	055045-09-5	74
5	Tricosane, 2-methyl-	338	C24H50	001928-30-9	70



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP120519\
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Misc :
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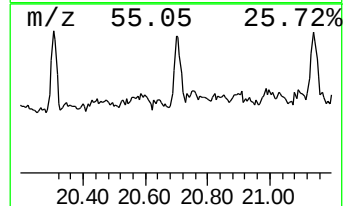
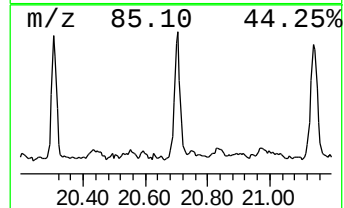
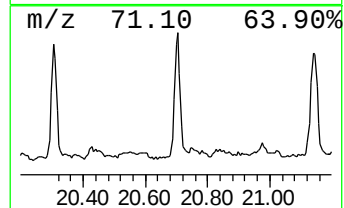
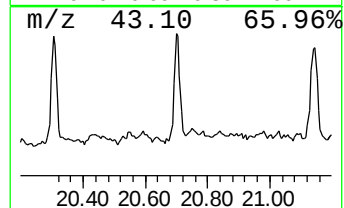
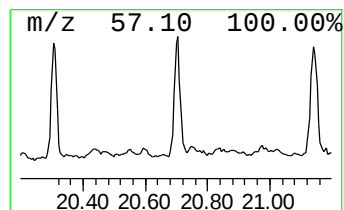
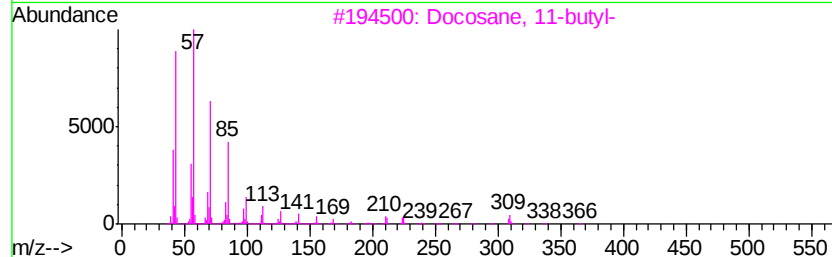
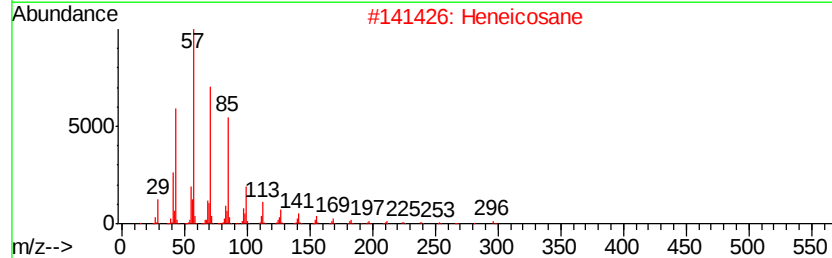
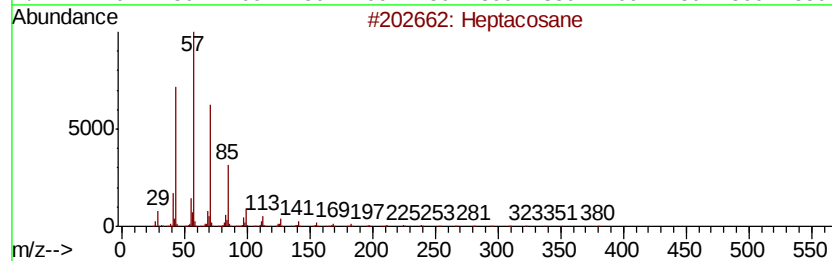
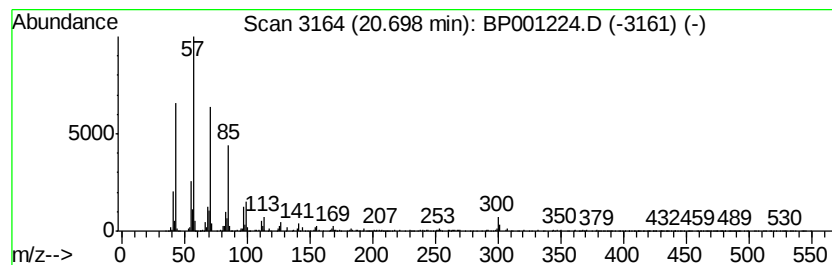
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ClientSampleId :
TH-7

Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP112719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Docosane, 11-butyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.70	6.96 ng	318980	Perylene-d12	21.02
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Heptacosane	380 C27H56	000593-49-7	96
2	Heneicosane	296 C21H44	000629-94-7	95
3	Docosane, 11-butyl-	366 C26H54	013475-76-8	93
4	Docosane, 9-butyl-	366 C26H54	055282-14-9	91
5	Heptadecane, 2,6,10,15-tetramethyl-	296 C21H44	054833-48-6	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP120519\
Data File : BP001224.D
Acq On : 05 Dec 2019 21:26
Operator : JU
Sample : K6120-01
Misc :
ALS Vial : 23 Sample Multiplier: 1

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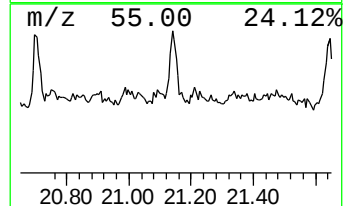
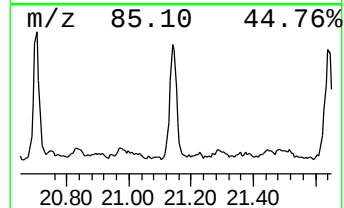
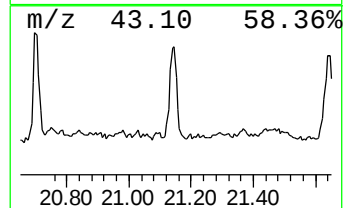
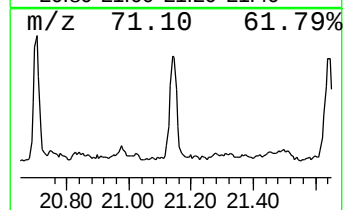
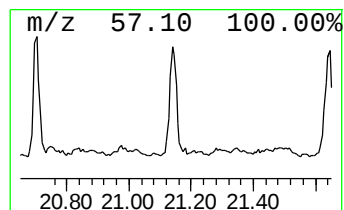
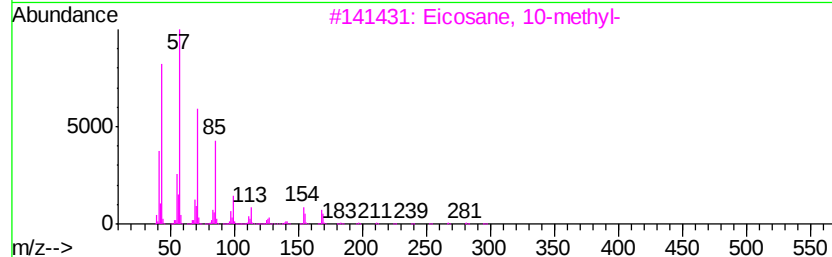
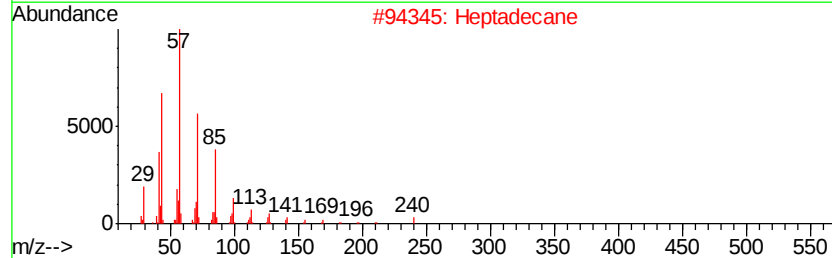
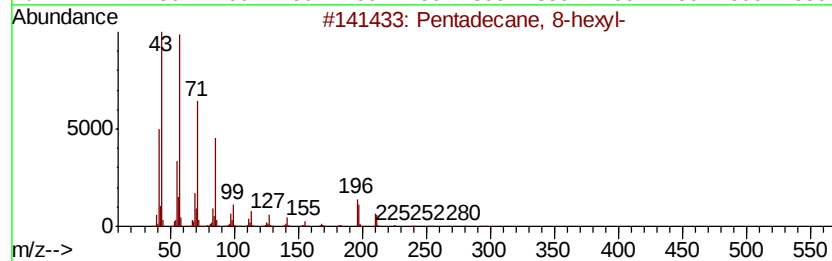
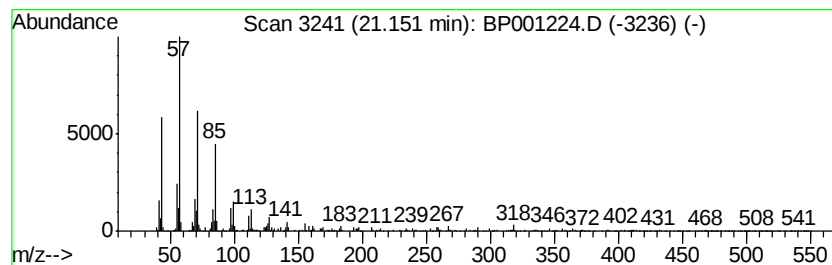
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Pentadecane, 8-hexyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.15	6.01 ng	275484	Perylene-d12	21.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	90
2		Heptadecane	240	C17H36	000629-78-7	90
3		Eicosane, 10-methyl-	296	C21H44	054833-23-7	90
4		Tetratetracontane	619	C44H90	007098-22-8	87
5		Tetradecane, 2,6,10-trimethyl-	240	C17H36	014905-56-7	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP120519\
Data File : BP001224.D
Acq On : 05 Dec 2019 21:26
Operator : JU
Sample : K6120-01
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
TH-7

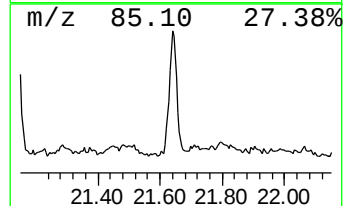
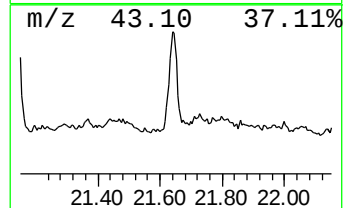
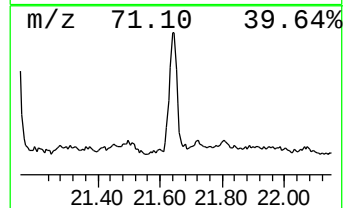
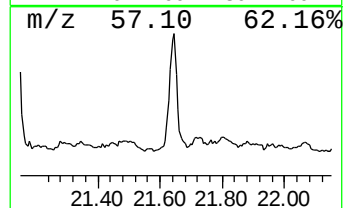
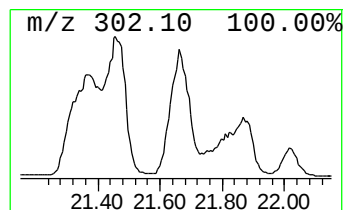
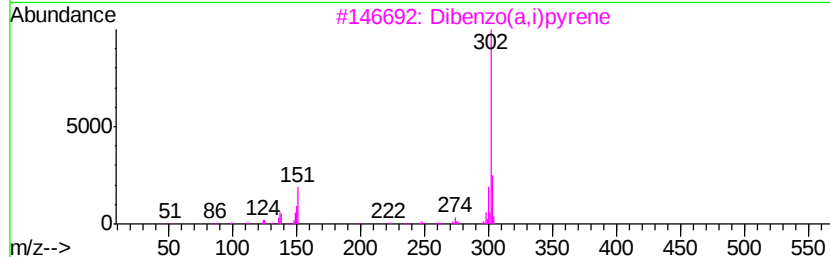
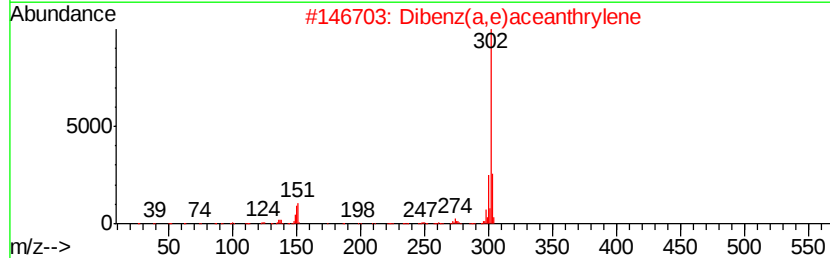
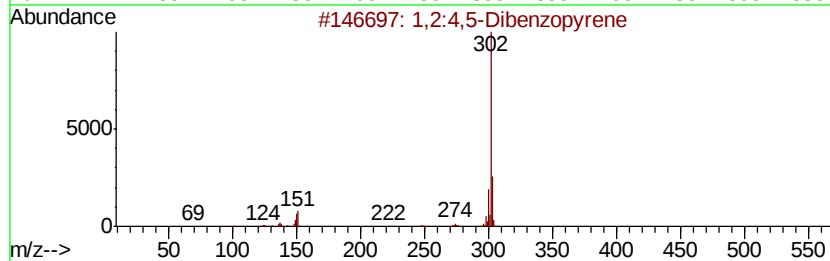
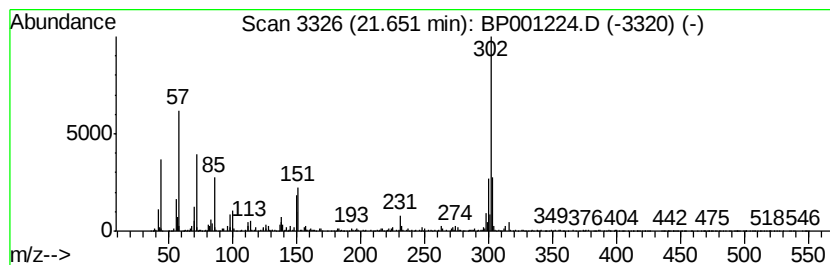
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 1,2:4,5-Dibenzopyrene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.65	17.74 ng	812656	Perylene-d12	21.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,2:4,5-Dibenzopyrene	302	C24H14	000192-65-4	93
2		Dibenz(a,e)aceanthrylene	302	C24H14	005385-75-1	89
3		Dibenzo(a,i)pyrene	302	C24H14	000189-55-9	81
4		3,4:8,9-Dibenzopyrene	302	C24H14	000189-64-0	74
5		Dibenzo[fg,op]naphthacene	302	C24H14	000192-51-8	68



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP120519\
Data File : BP001224.D
Acq On : 05 Dec 2019 21:26
Operator : JU
Sample : K6120-01
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
TH-7

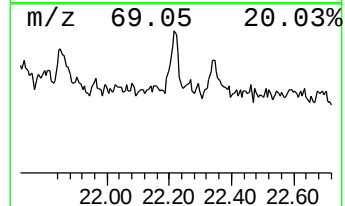
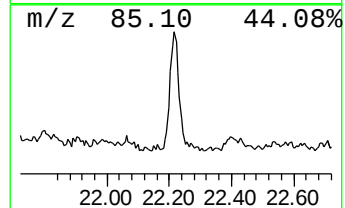
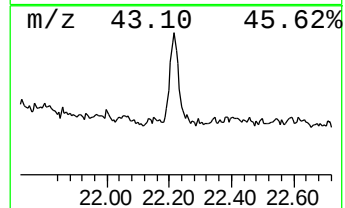
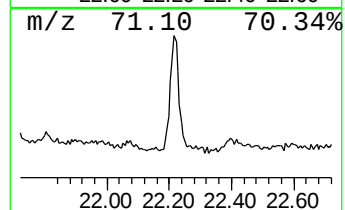
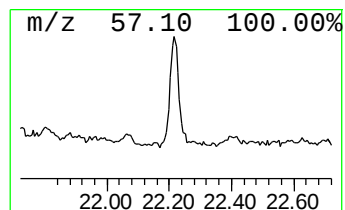
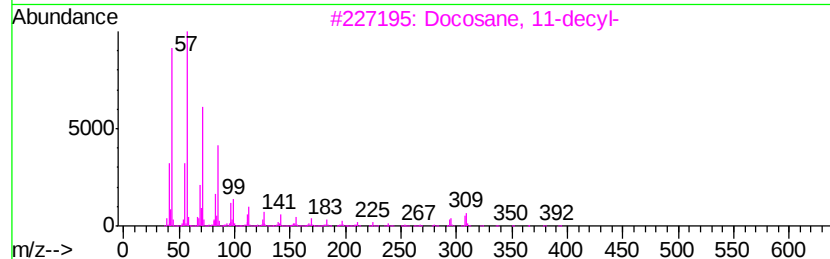
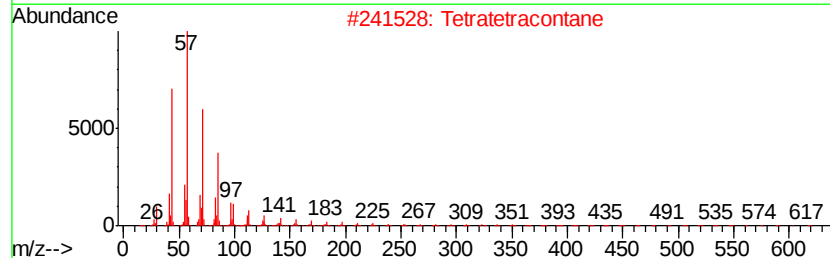
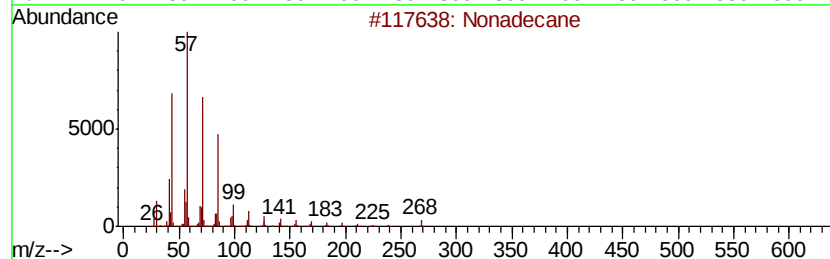
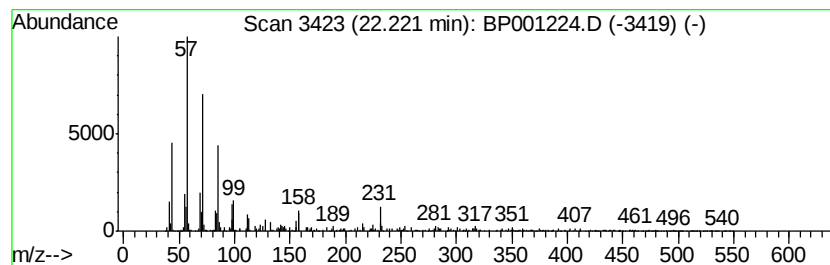
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 Tetratetracontane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.22	4.50 ng	206065	Perylene-d12	21.02

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonadecane	268	C19H40	000629-92-5	94
2			Tetratetracontane	619	C44H90	007098-22-8	86
3			Docosane, 11-decyl-	451	C32H66	055401-55-3	86
4			Docosane	310	C22H46	000629-97-0	81
5			Docosane, 1,22-dibromo-	466	C22H44Br2	034540-49-3	74



Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP120519\
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 Acq On : 05 Dec 2019 21:26
 Operator : JU
 Sample : K6120-01
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 TH-7

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP112719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.63	17.1 ng		475350	1	8.31	555645	20.0
unknown7.95	7.95	86.4 ng		2400920	1	8.31	555645	20.0
n-Hexadecanoic acid	16.49	4.2 ng		176182	4	15.60	830231	20.0
Benzene, 1,1'-(1,...	17.35	2.1 ng		87358	4	15.60	830231	20.0
Heneicosane	18.26	2.1 ng		85069	5	19.25	818020	20.0
Heptacosane	18.71	3.2 ng		129803	5	19.25	818020	20.0
Octadecyl trifluo...	19.13	15.3 ng		626287	5	19.25	818020	20.0
Nonadecane	19.54	4.0 ng		162417	5	19.25	818020	20.0
Octadecane, 5,14-...	19.93	6.8 ng		278308	5	19.25	818020	20.0
unknown20.12	20.12	5.9 ng		240306	5	19.25	818020	20.0
Tetracosane	20.30	3.5 ng		161674	6	21.02	916357	20.0
Docosane, 11-butyl-	20.70	7.0 ng		318980	6	21.02	916357	20.0
Pentadecane, 8-he...	21.15	6.0 ng		275484	6	21.02	916357	20.0
1,2:4,5-Dibenzopy...	21.65	17.7 ng		812656	6	21.02	916357	20.0
Tetratetracontane	22.22	4.5 ng		206065	6	21.02	916357	20.0