

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP011620\
 Data File : BP001633.D
 Acq On : 16 Jan 2020 18:09
 Operator : JU
 Sample : L1145-01
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 MH-5

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

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

Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP010820.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	2.870	9	14	22	rBV	55585	100510	4.05%	0.571%
2	2.946	22	27	41	rVB	580334	766387	30.90%	4.352%
3	4.793	337	341	349	rVB	55618	79561	3.21%	0.452%
4	5.258	414	420	430	rVB	400176	575594	23.21%	3.269%
5	6.822	680	686	696	rBV	252019	421984	17.02%	2.396%
6	7.169	738	745	758	rBV	725697	1160944	46.81%	6.593%
7	7.628	815	823	832	rBV	196984	310591	12.52%	1.764%
8	7.946	870	877	887	rVB	614912	986702	39.79%	5.604%
9	8.775	1006	1018	1033	rBV	509132	865950	34.92%	4.918%
10	10.404	1288	1295	1304	rBV	242722	400465	16.15%	2.274%
11	12.893	1711	1718	1730	rBV	1134841	1717311	69.25%	9.753%
12	13.198	1762	1770	1778	rBV	39035	64294	2.59%	0.365%
13	13.740	1856	1862	1874	rBV	76356	116780	4.71%	0.663%
14	14.275	1946	1953	1961	rBV	368915	553348	22.31%	3.143%
15	15.057	2081	2086	2091	rVB	18458	25212	1.02%	0.143%
16	15.781	2201	2209	2220	rBV	1033378	1472274	59.37%	8.361%
17	16.598	2344	2348	2353	rBV	39493	54860	2.21%	0.312%
18	17.039	2417	2423	2434	rBV	487515	662713	26.72%	3.764%
19	17.551	2505	2510	2514	rBV2	31844	49859	2.01%	0.283%
20	17.845	2556	2560	2567	rBV6	18468	31911	1.29%	0.181%
21	17.975	2577	2582	2596	rBV2	120179	204080	8.23%	1.159%
22	19.627	2857	2863	2874	rBV	2105279	2479888	100.00%	14.084%
23	19.957	2916	2919	2922	rVB	25296	26057	1.05%	0.148%
24	20.439	2998	3001	3005	rVB	46901	49589	2.00%	0.282%
25	20.892	3073	3078	3083	rBV	312489	441019	17.78%	2.505%
26	21.145	3116	3121	3128	rBV	479123	639735	25.80%	3.633%
27	21.322	3147	3151	3164	rVB	184941	256819	10.36%	1.459%
28	21.757	3221	3225	3235	rBV	210442	290410	11.71%	1.649%
29	22.227	3301	3305	3311	rVB	290717	373698	15.07%	2.122%
30	22.357	3322	3327	3333	rVB	312464	427989	17.26%	2.431%
31	22.739	3387	3392	3398	rVB	263685	379476	15.30%	2.155%
32	23.316	3484	3490	3494	rBV	255335	407410	16.43%	2.314%
33	23.368	3494	3499	3511	rVB	320397	597684	24.10%	3.394%
34	23.974	3596	3602	3610	rVB	193845	356006	14.36%	2.022%

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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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35	24.057	3612	3616	3625	rVB7	31515	79170	3.19%	0.450%
36	24.733	3726	3731	3738	rVB	92516	182120	7.34%	1.034%

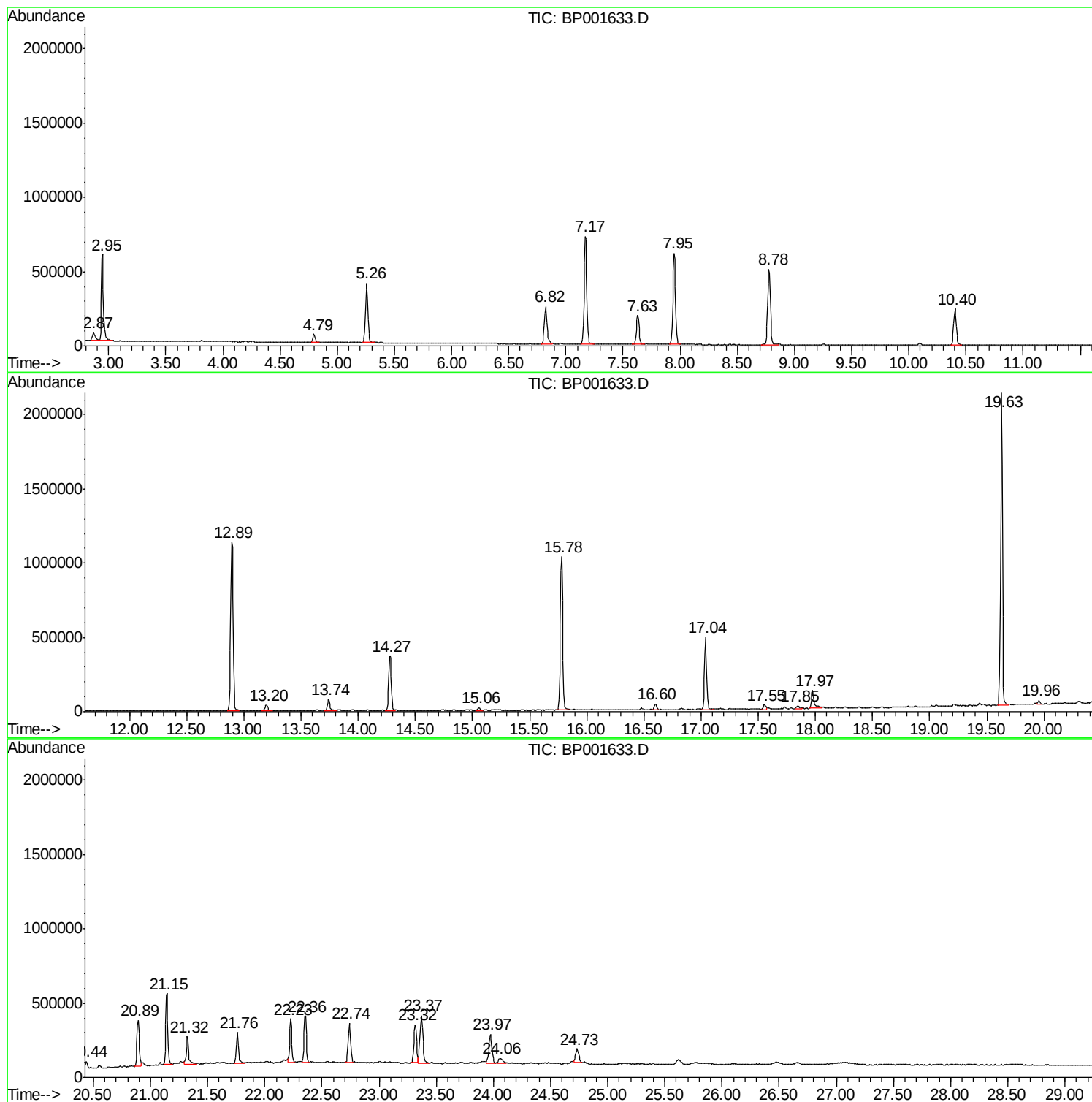
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP010820.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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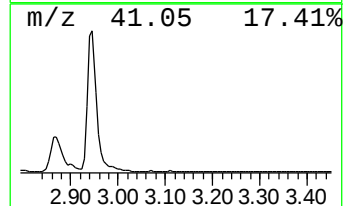
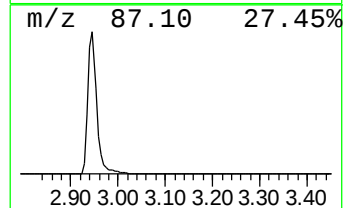
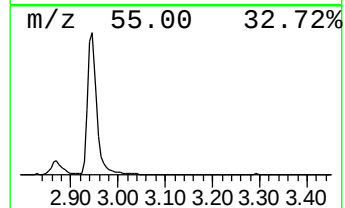
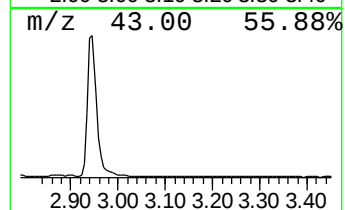
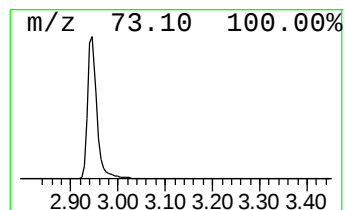
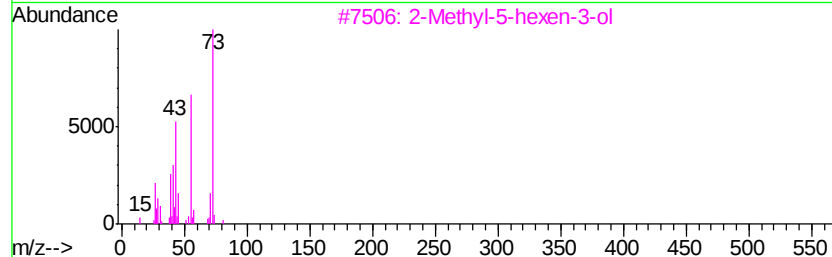
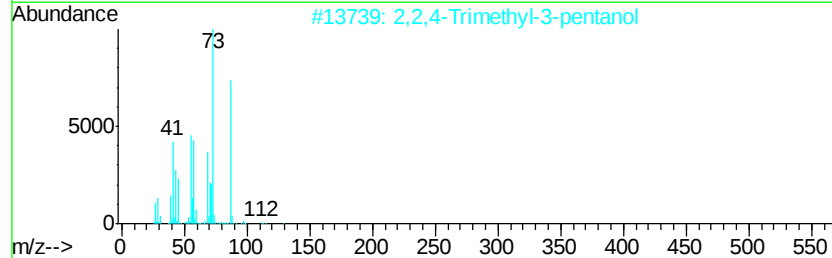
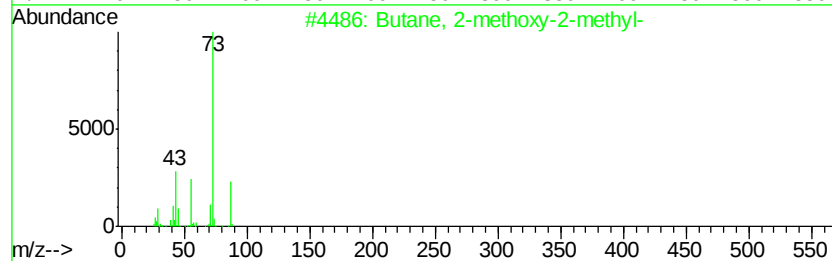
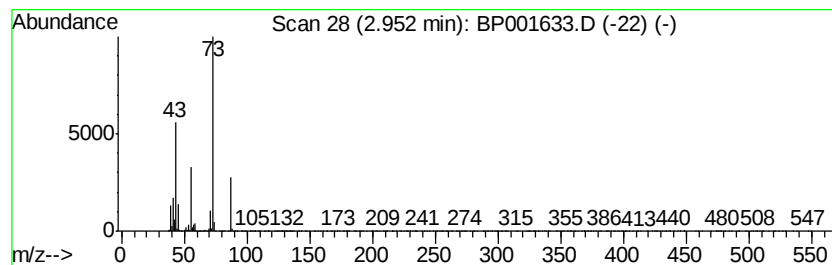
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.95	49.35 ng	766387	1,4-Dichlorobenzene-d4	7.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
3		2-Methyl-5-hexen-3-ol	114	C7H14O	032815-70-6	23
4		Silane, tetramethyl-	88	C4H12Si	000075-76-3	23
5		Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9



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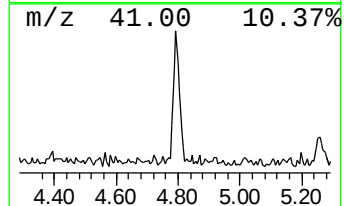
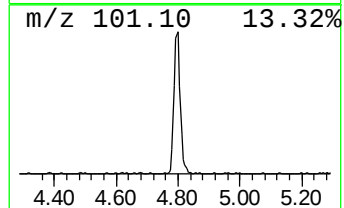
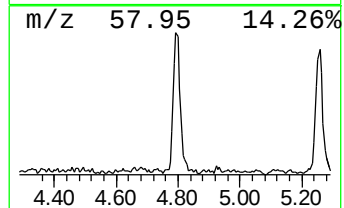
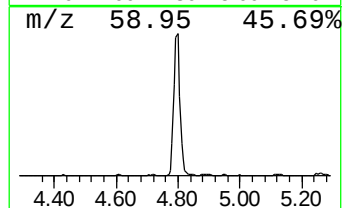
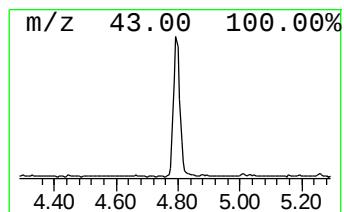
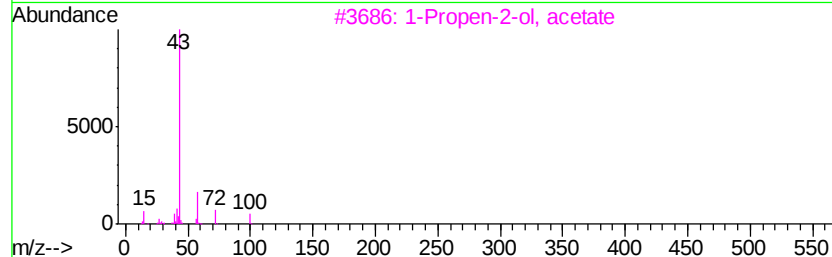
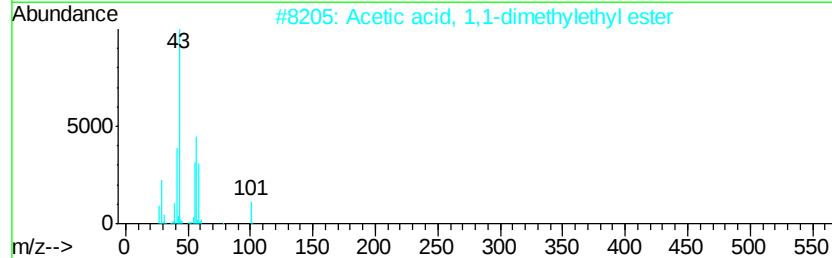
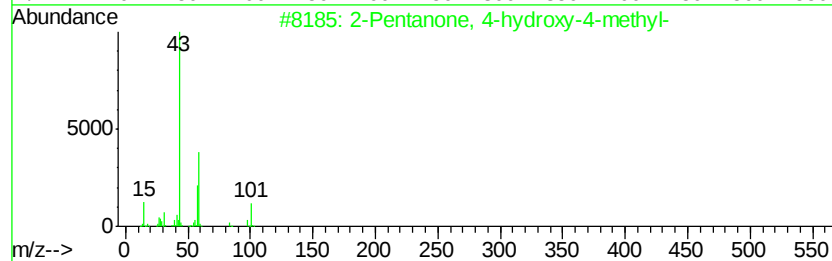
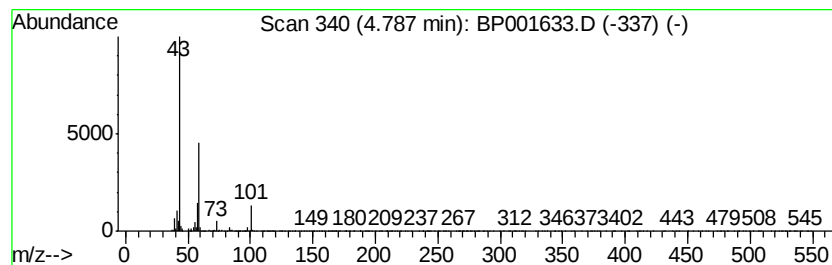
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.79	5.12 ng	79561	1,4-Dichlorobenzene-d4	7.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	28
3		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
4		(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	9
5		Propanenitrile, 3-ethoxy-	99	C5H9NO	002141-62-0	9



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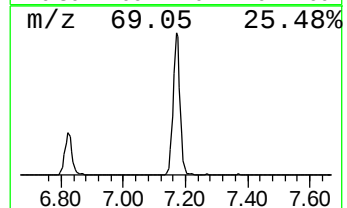
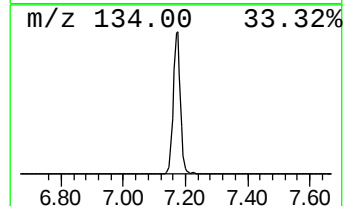
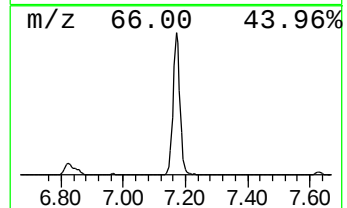
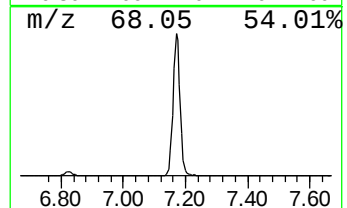
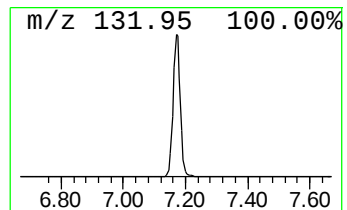
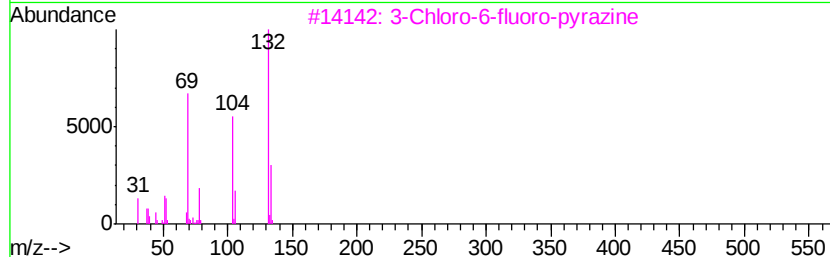
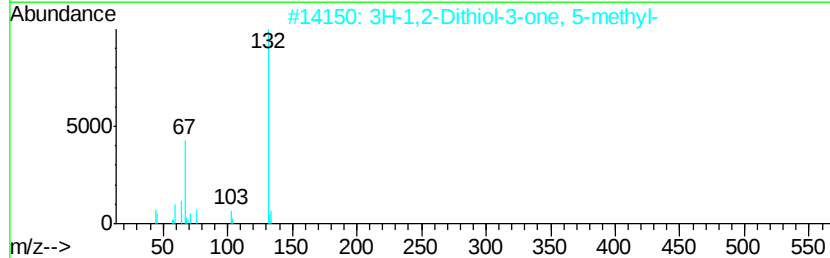
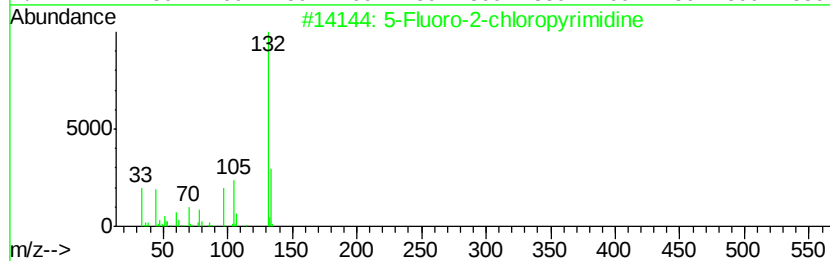
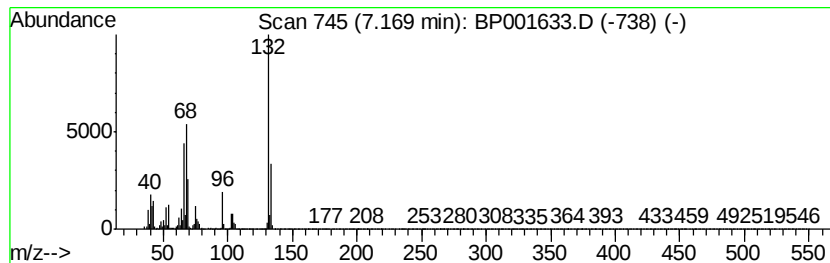
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 unknown7.17 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.17	74.76 ng	1160940	1,4-Dichlorobenzene-d4	7.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	14
2		3H-1,2-Dithiol-3-one, 5-methyl-	132	C4H4OS2	003620-08-4	14
3		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	14
4		1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	12
5		1,3,4-Thiadiazole-2(3H)-thione, ...	132	C3H4N2S2	029490-19-5	10



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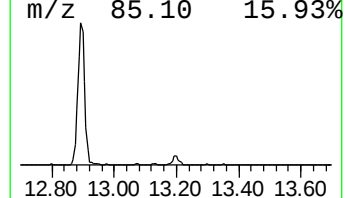
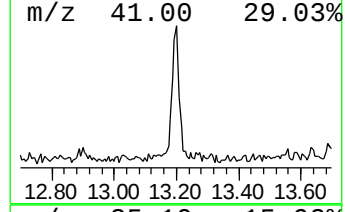
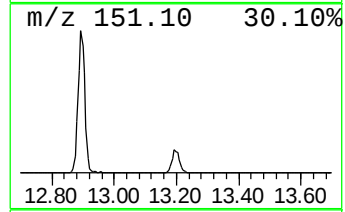
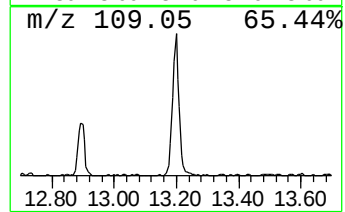
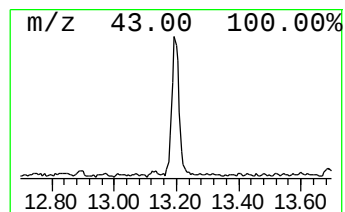
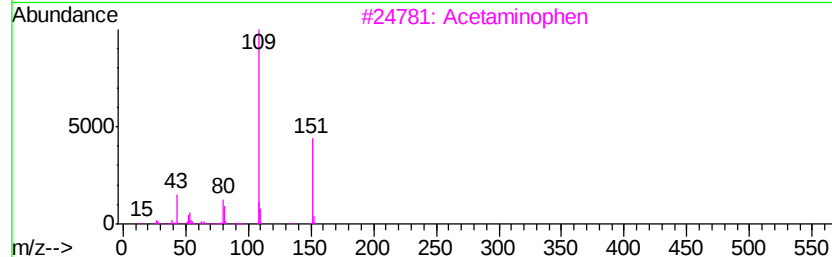
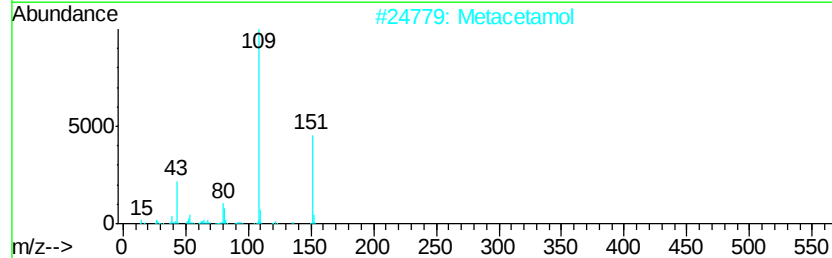
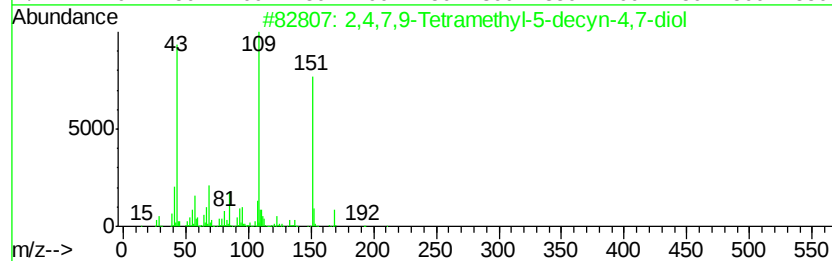
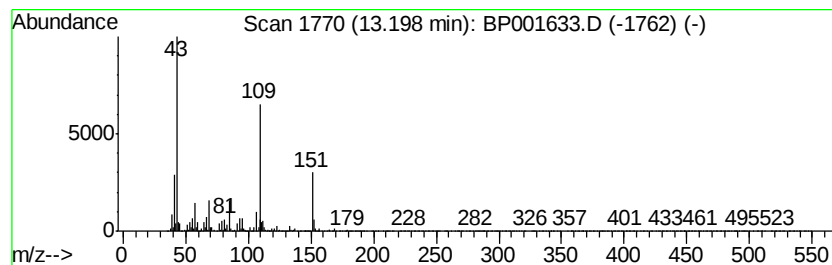
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 2,4,7,9-Tetramethyl-5-decyn... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.20	2.32 ng	64294	Acenaphthene-d10	14.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2,4,7,9-Tetramethyl-5-decyn-4,7-...	226	C14H26O2	000126-86-3	86
2		Metacetamol	151	C8H9NO2	000621-42-1	52
3		Acetaminophen	151	C8H9NO2	000103-90-2	47
4		1H-Pyrazole, 4-[4-(3-chloropheno...	278	C15H19ClN2O	1000302-33-9	40
5		m-Isopropoxyaniline	151	C9H13NO	041406-00-2	40



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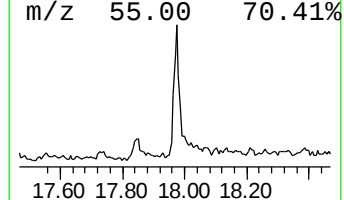
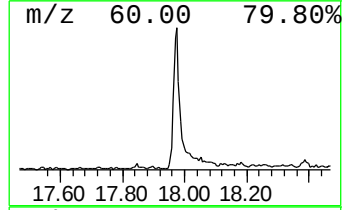
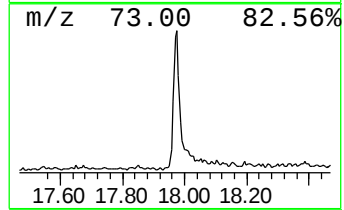
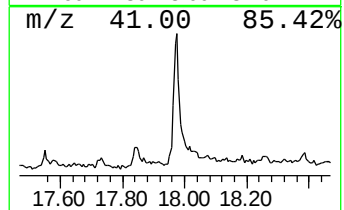
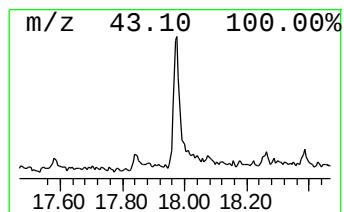
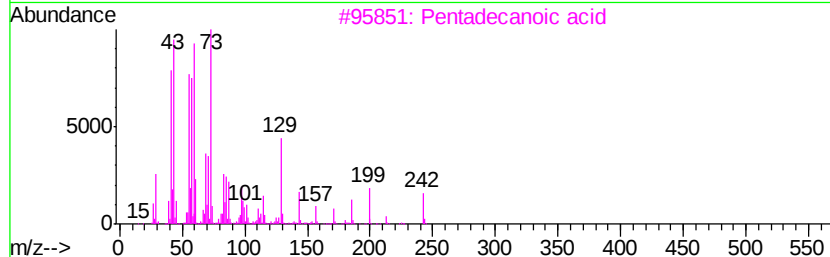
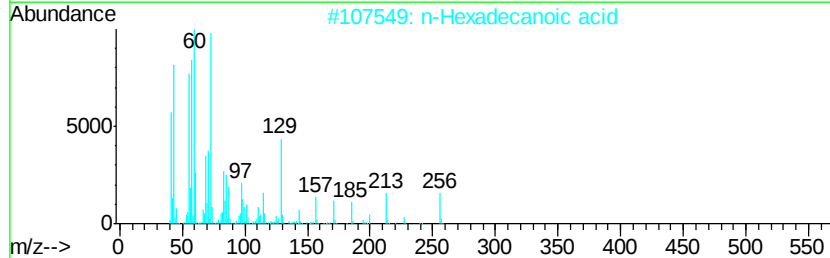
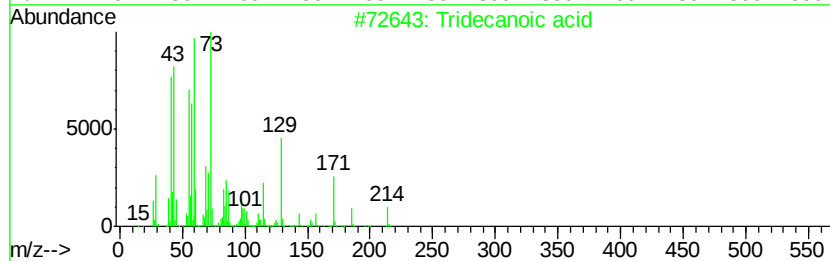
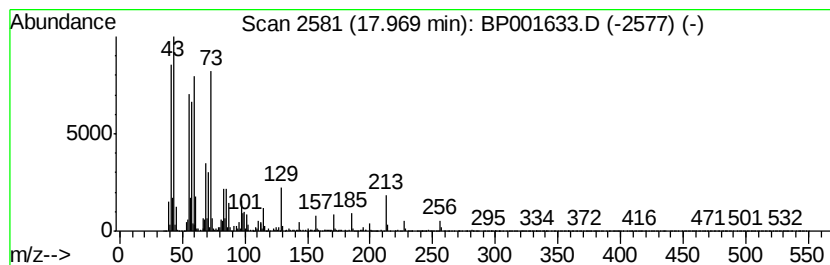
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 7 Tridecanoic acid Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.97	6.16 ng	204080	Phenanthrene-d10	17.04

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tridecanoic acid	214	C13H26O2	000638-53-9	95
2		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	91
3		Pentadecanoic acid	242	C15H30O2	001002-84-2	74
4		Undecanoic acid	186	C11H22O2	000112-37-8	70
5		Tetradecanoic acid	228	C14H28O2	000544-63-8	64



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Sample : L1145-01
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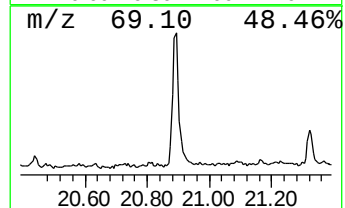
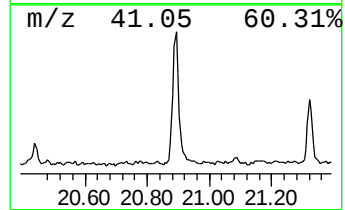
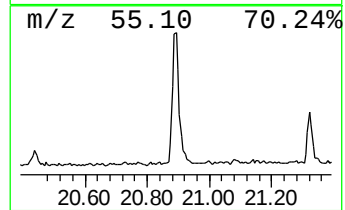
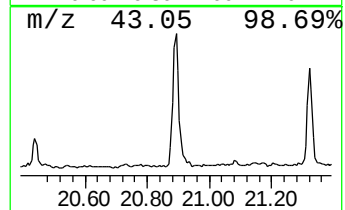
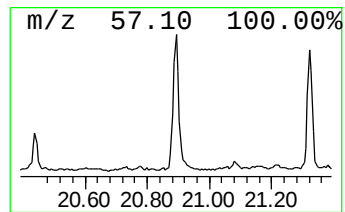
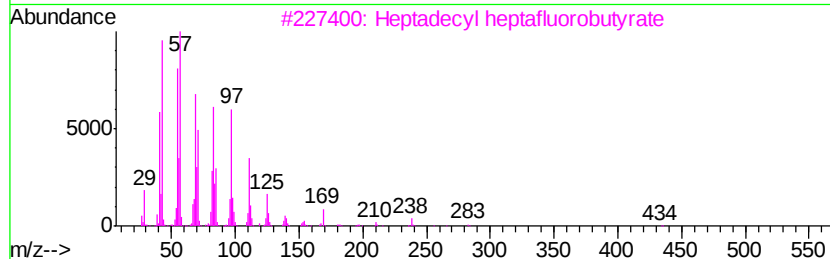
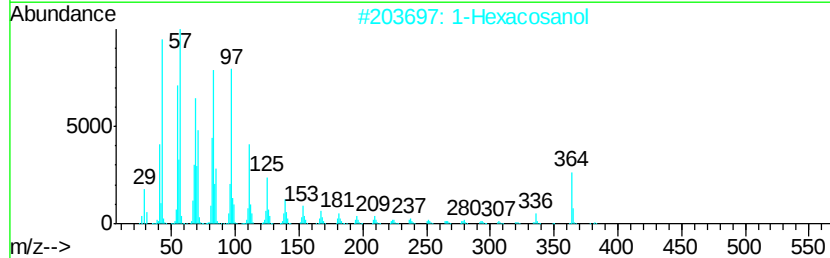
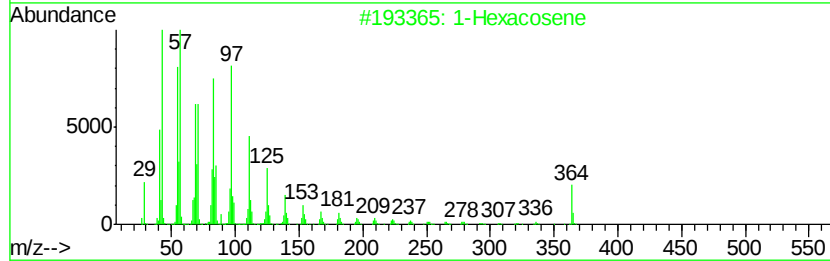
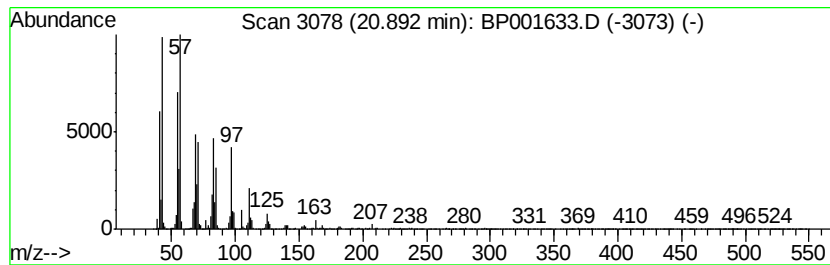
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP010820.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 1-Hexacosene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.89	13.79 ng	441019	Chrysene-d12	21.14
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1-Hexacosene	364 C26H52	018835-33-1	95
2	1-Hexacosanol	382 C26H54O	000506-52-5	91
3	Heptadecyl heptafluorobutvrate	452 C21H35F7O2	959085-66-6	91
4	Cyclohexadecane, 1,2-diethyl-	280 C20H40	1000155-85-3	89
5	E-15-Heptadecenal	252 C17H32O	1000130-97-9	81



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP011620\
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Acq On : 16 Jan 2020 18:09
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ClientSampleId :
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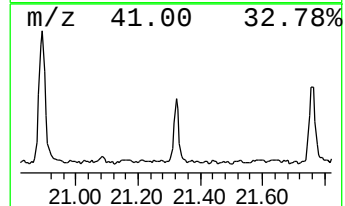
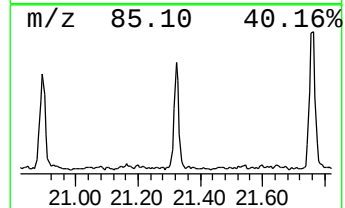
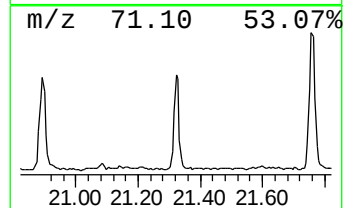
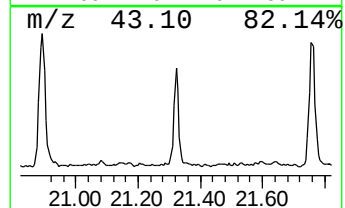
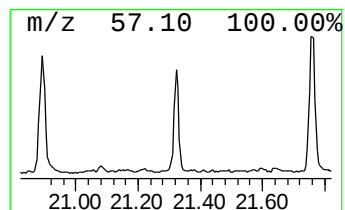
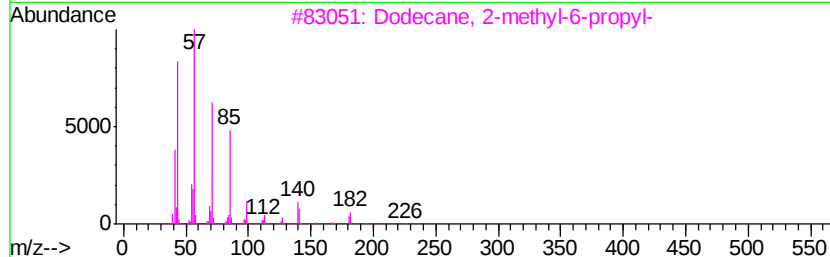
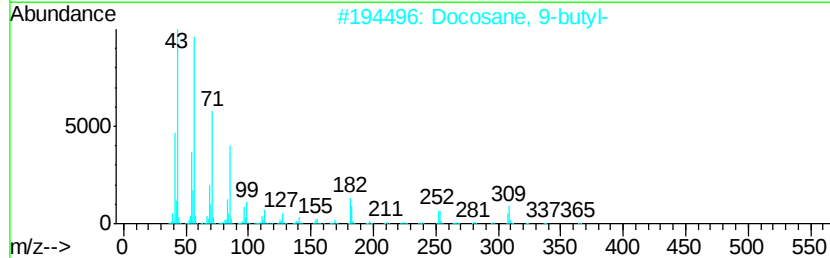
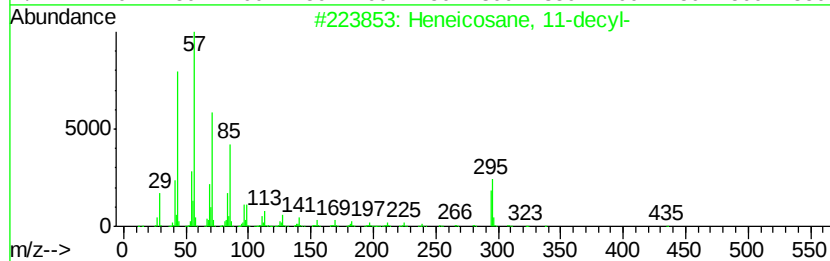
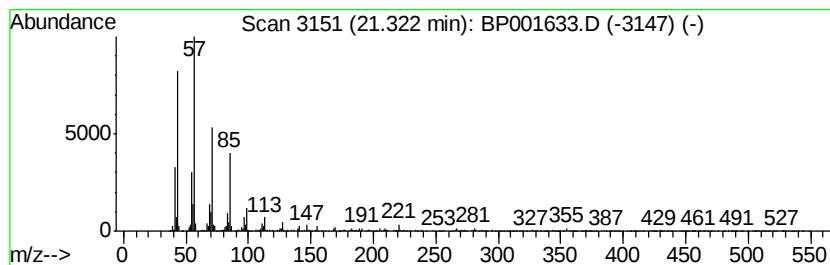
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Heneicosane, 11-decyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.32	8.03 ng	256819	Chrysene-d12	21.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane, 11-decyl-	437	C31H64	055320-06-4	93
2		Docosane, 9-butyl-	366	C26H54	055282-14-9	91
3		Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	91
4		Pentacosane	352	C25H52	000629-99-2	90
5		Docosane, 5-butyl-	366	C26H54	055282-16-1	89



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP011620\
Data File : BP001633.D
Acq On : 16 Jan 2020 18:09
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Misc :
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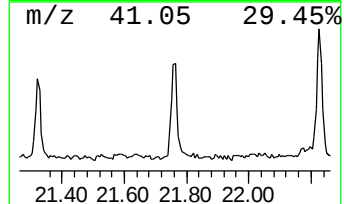
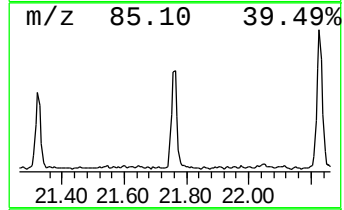
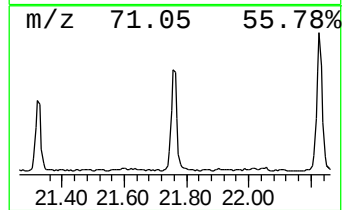
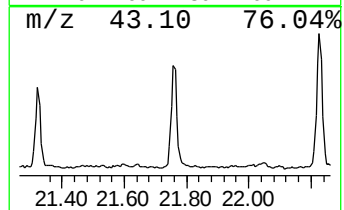
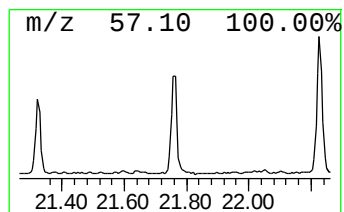
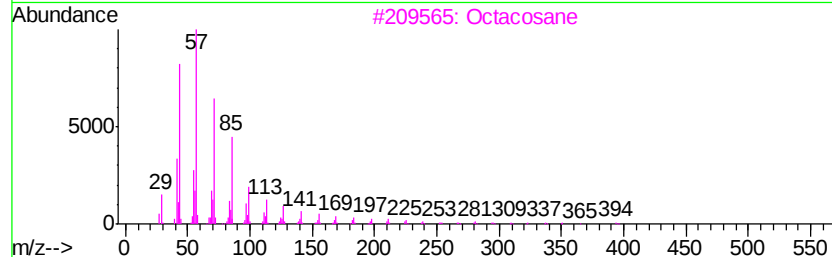
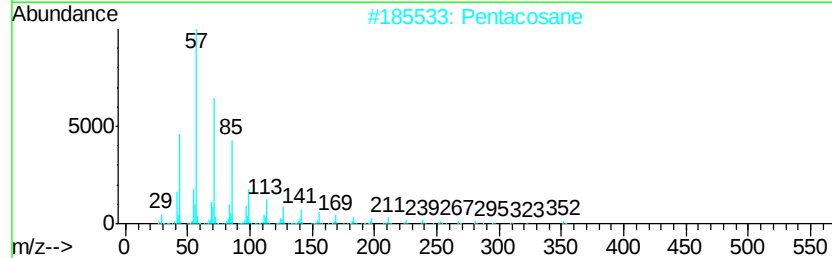
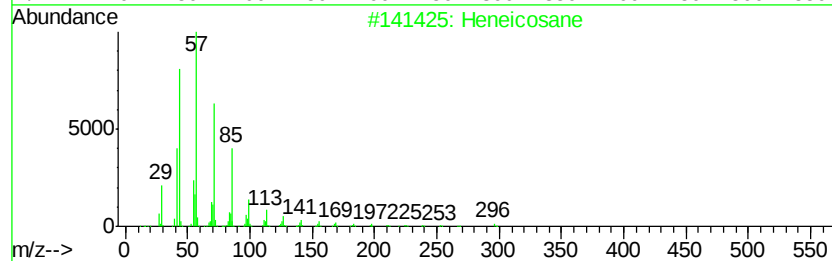
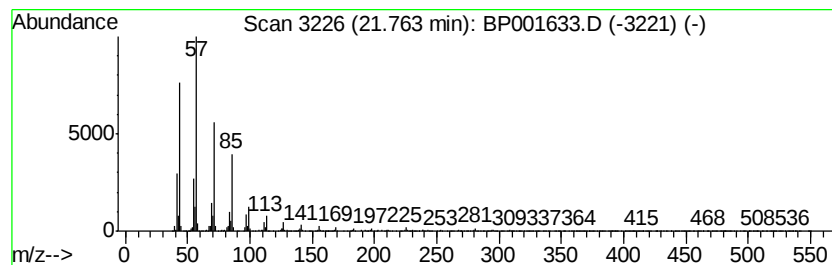
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Heneicosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.76	9.08 ng	290410	Chrysene-d12	21.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	98
2		Pentacosane	352	C25H52	000629-99-2	97
3		Octacosane	394	C28H58	000630-02-4	95
4		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	95
5		Eicosane, 10-methyl-	296	C21H44	054833-23-7	95



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP011620\
Data File : BP001633.D
Acq On : 16 Jan 2020 18:09
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Misc :
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Instrument :
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ClientSampleId :
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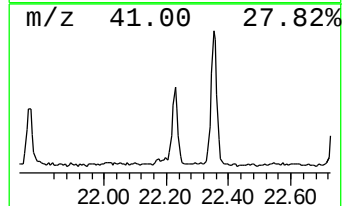
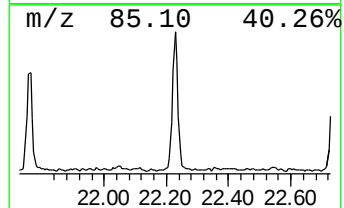
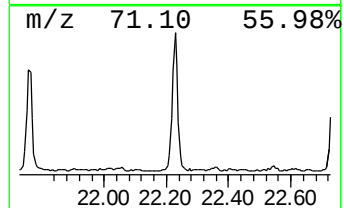
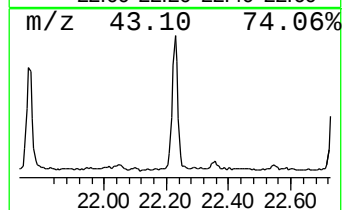
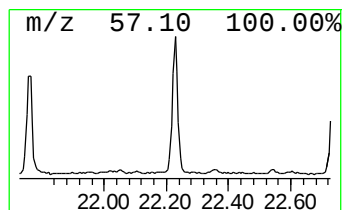
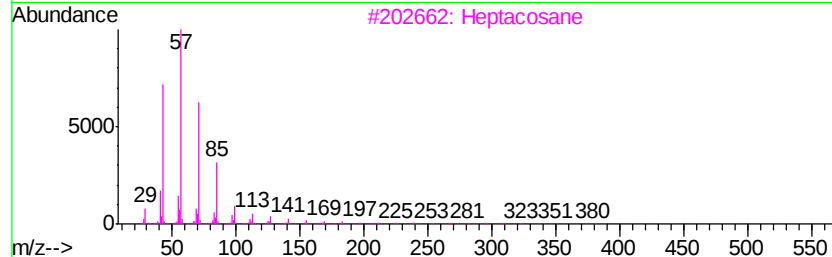
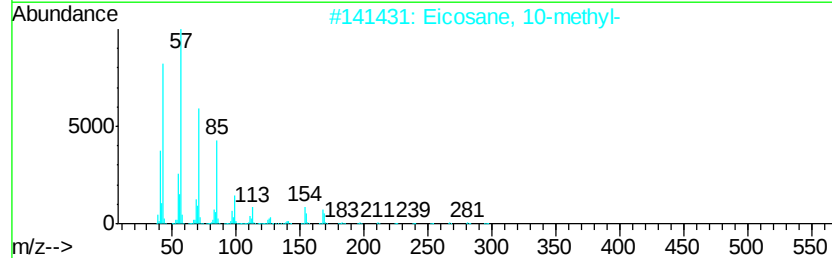
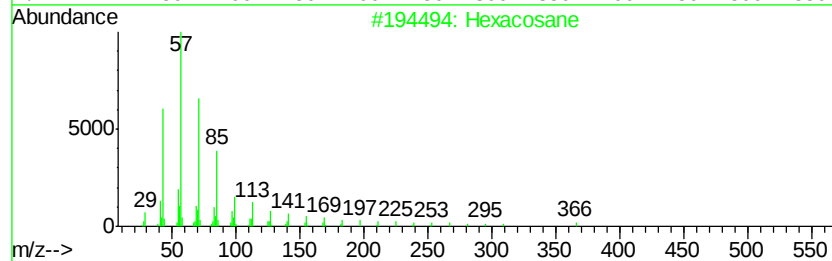
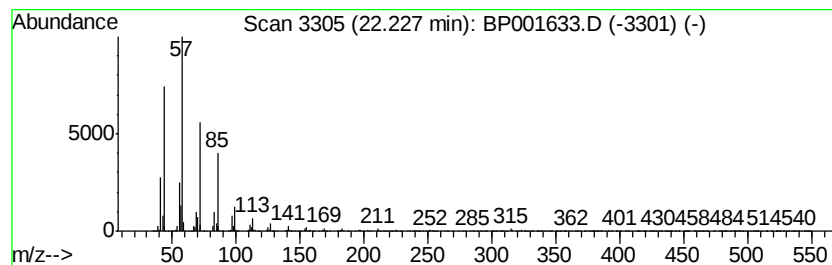
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Hexacosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.23	11.68 ng	373698	Chrysene-d12	21.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexacosane	366	C26H54	000630-01-3	95
2		Eicosane, 10-methyl-	296	C21H44	054833-23-7	94
3		Heptacosane	380	C27H56	000593-49-7	91
4		Tetracosane	338	C24H50	000646-31-1	91
5		Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP011620\
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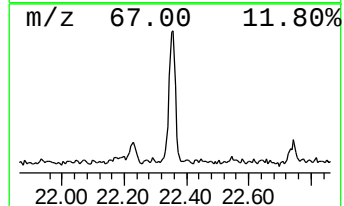
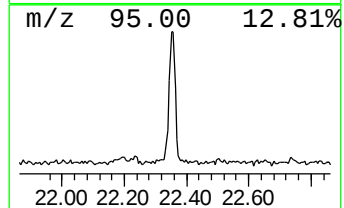
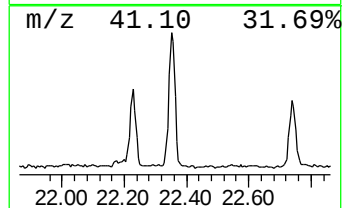
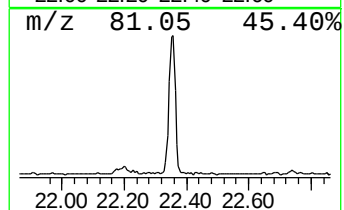
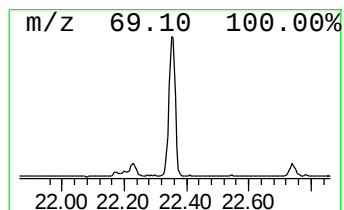
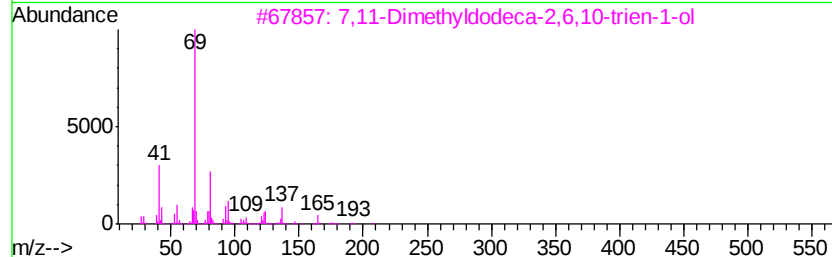
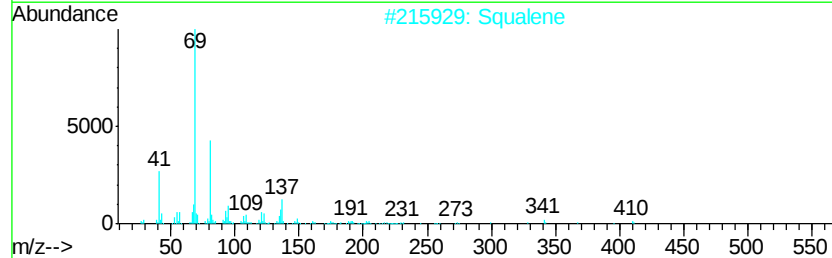
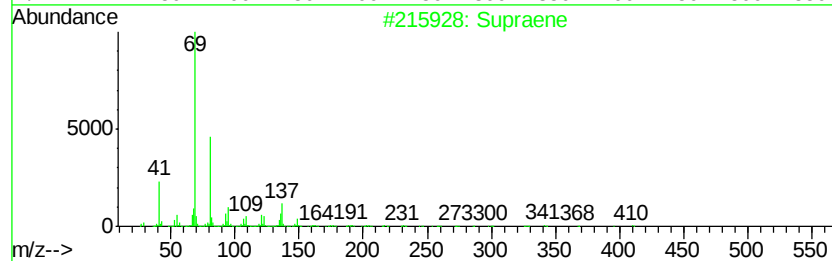
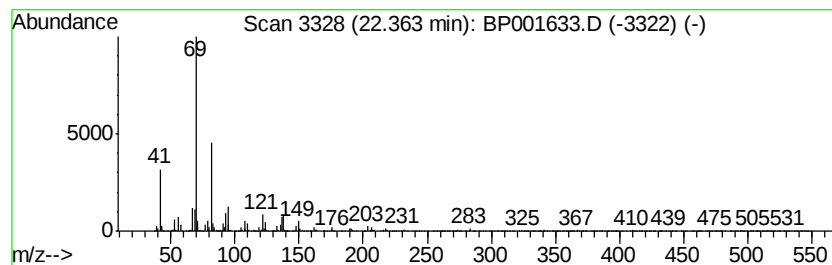
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 Supraene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.36	14.32 ng	427989	Perylene-d12	23.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Supraene	410	C30H50	007683-64-9	91
2		Squalene	410	C30H50	000111-02-4	90
3		7,11-Dimethyldodeca-2,6,10-trien...	208	C14H24O	125529-10-4	72
4		3,7,11-Tridecatrienitrile, 4,8...	231	C16H25N	006006-01-5	64
5		2,3,5-Trimethyl-2,3,5-hexanetric...	203	C12H17N3	003974-79-6	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP011620\
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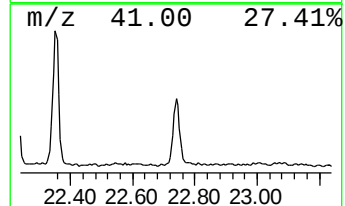
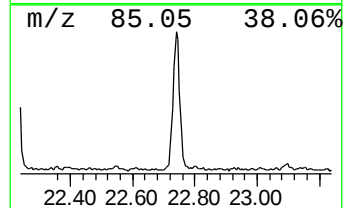
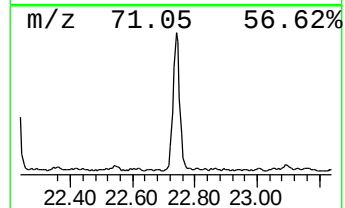
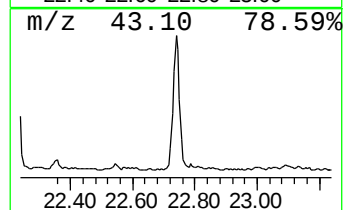
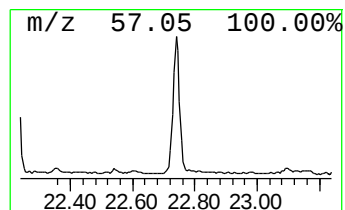
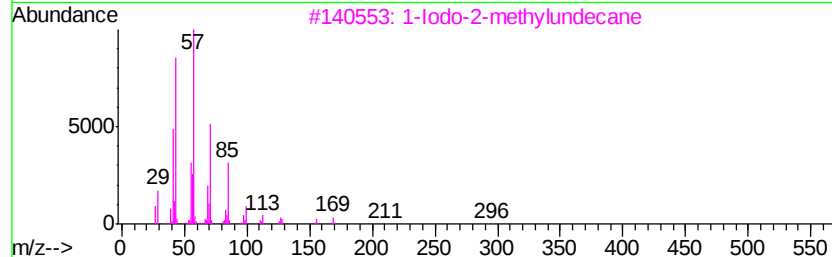
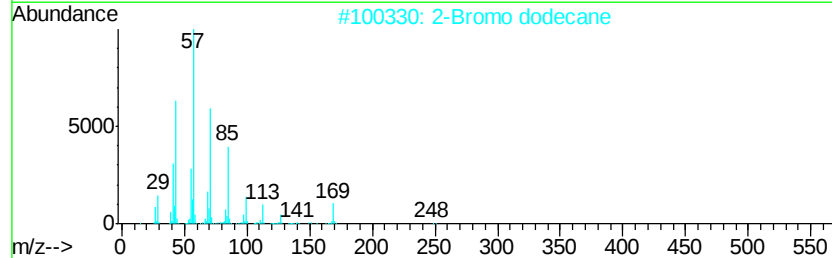
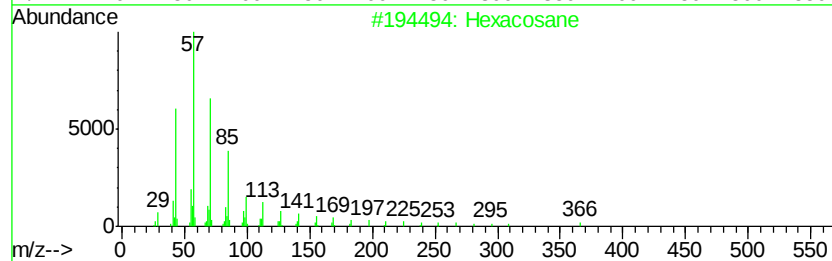
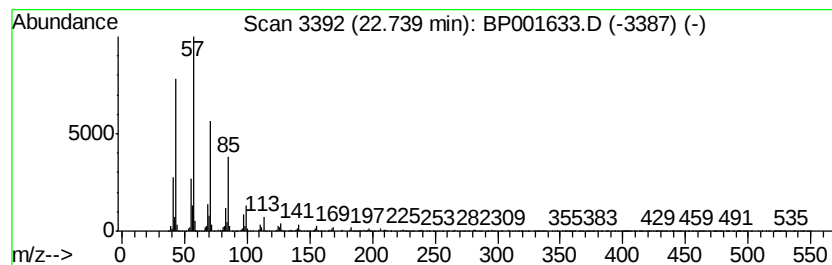
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 2-Bromo dodecane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.74	12.70 ng	379476	Perylene-d12	23.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexacosane	366	C26H54	000630-01-3	94
2		2-Bromo dodecane	248	C12H25Br	013187-99-0	94
3		1-Iodo-2-methylundecane	296	C12H25I	073105-67-6	94
4		Docosane, 9-butyl-	366	C26H54	055282-14-9	94
5		Octadecane, 5,14-dibutyl-	366	C26H54	055282-13-8	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP011620\
Data File : BP001633.D
Acq On : 16 Jan 2020 18:09
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ClientSampleId :
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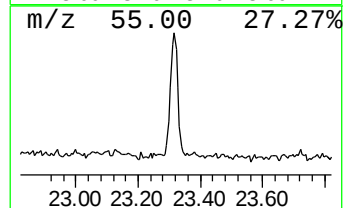
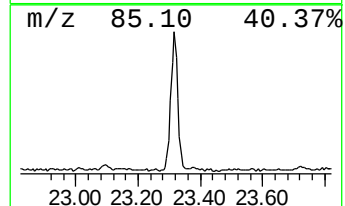
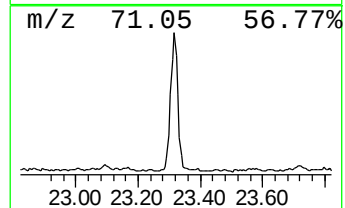
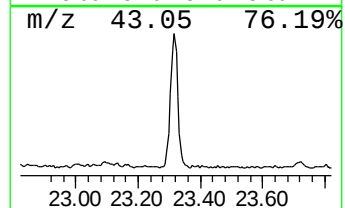
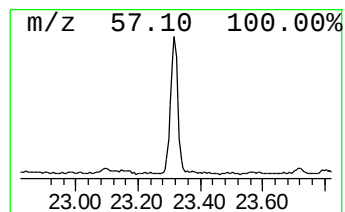
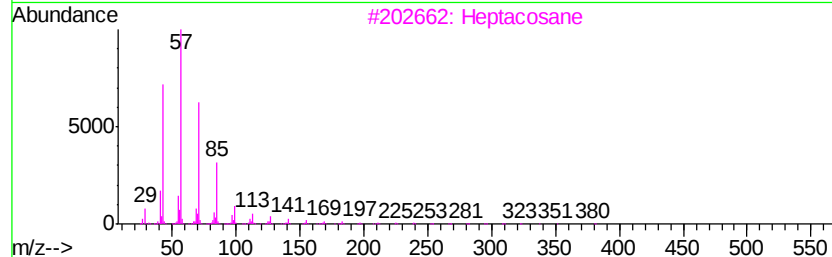
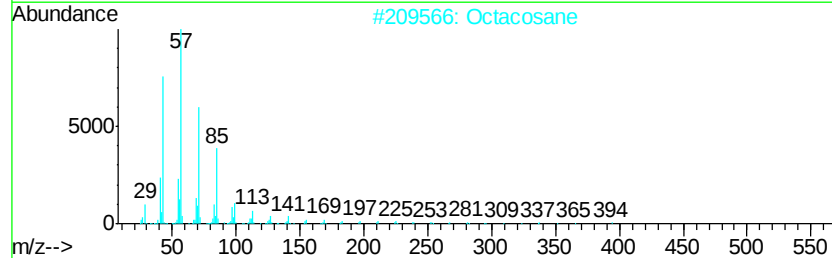
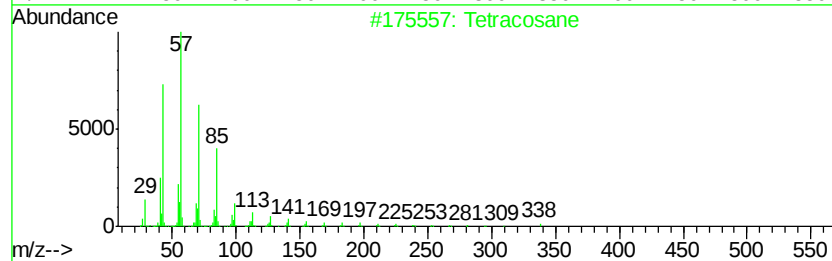
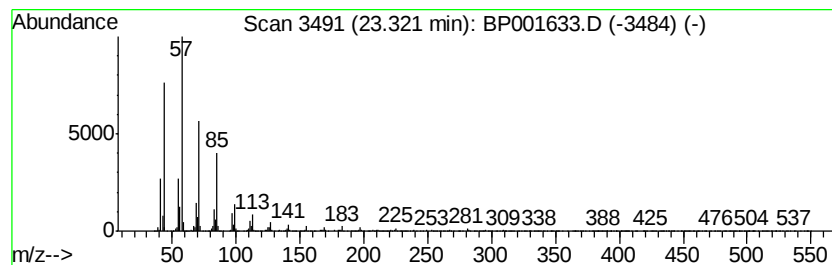
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 2-methyloctacosane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.32	13.63 ng	407410	Perylene-d12	23.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetracosane	338	C24H50	000646-31-1	91
2		Octacosane	394	C28H58	000630-02-4	91
3		Heptacosane	380	C27H56	000593-49-7	91
4		2-methyloctacosane	408	C29H60	1000376-72-8	91
5		Hexacosane	366	C26H54	000630-01-3	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP011620\
 Data File : BP001633.D
 Acq On : 16 Jan 2020 18:09
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 Sample : L1145-01
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
 MH-5

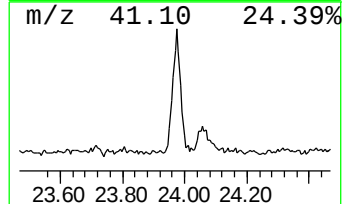
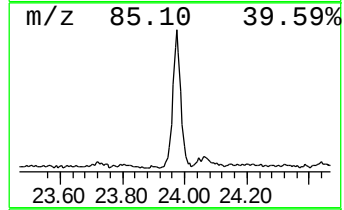
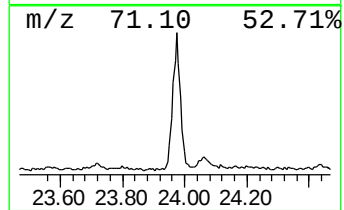
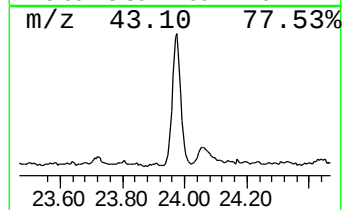
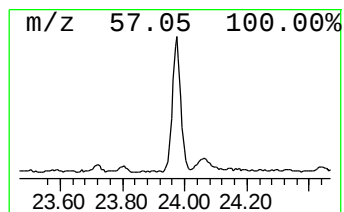
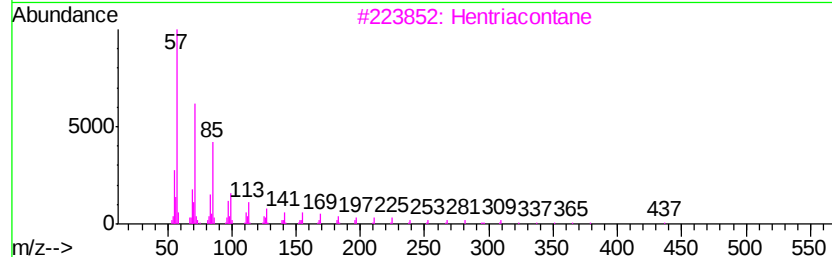
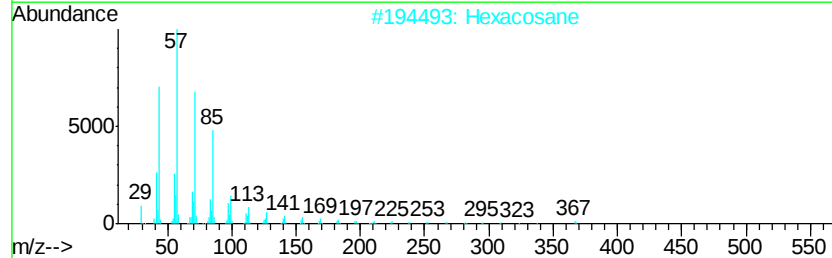
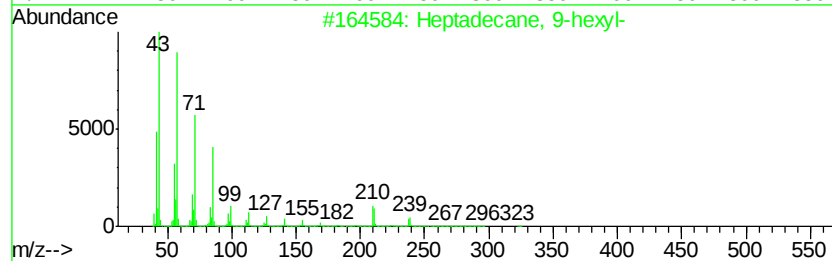
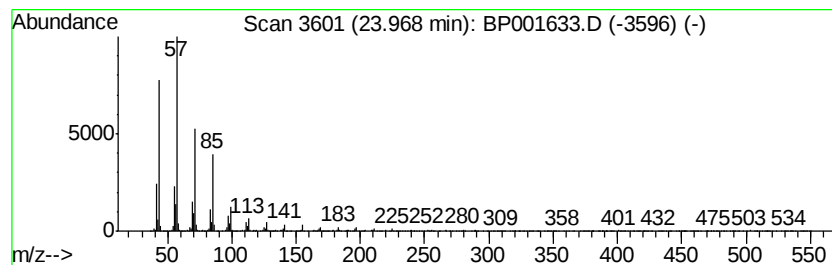
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 15 Eicosane, 7-hexyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.97	11.91 ng	356006	Perylene-d12	23.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane, 9-hexyl-	324	C23H48	055124-79-3	93
2		Hexacosane	366	C26H54	000630-01-3	91
3		Hentriacontane	437	C31H64	000630-04-6	91
4		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	90
5		Eicosane, 7-hexyl-	366	C26H54	055333-99-8	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP011620\
Data File : BP001633.D
Acq On : 16 Jan 2020 18:09
Operator : JU
Sample : L1145-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
MH-5

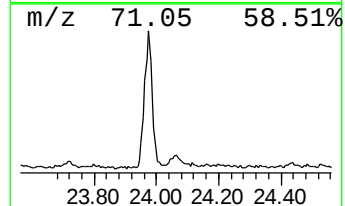
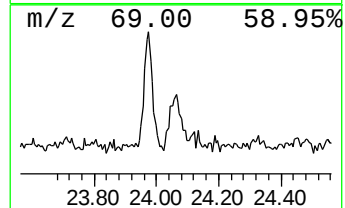
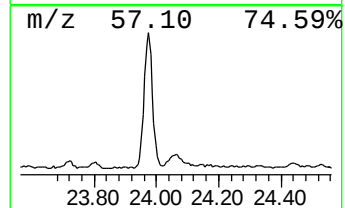
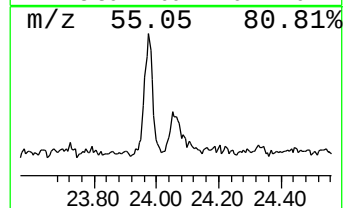
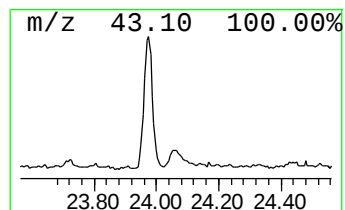
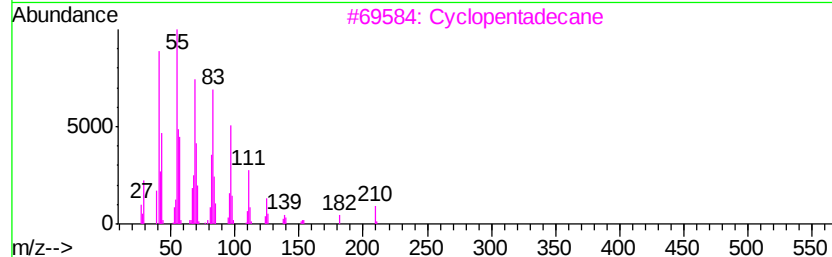
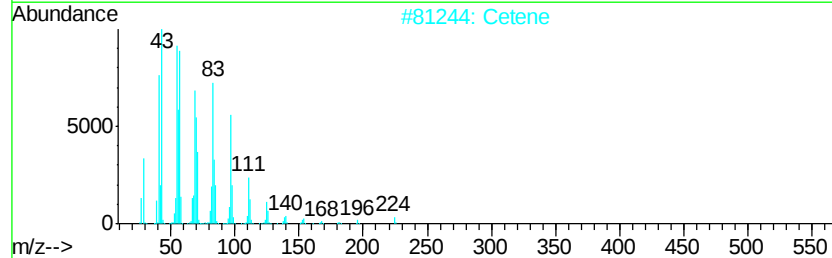
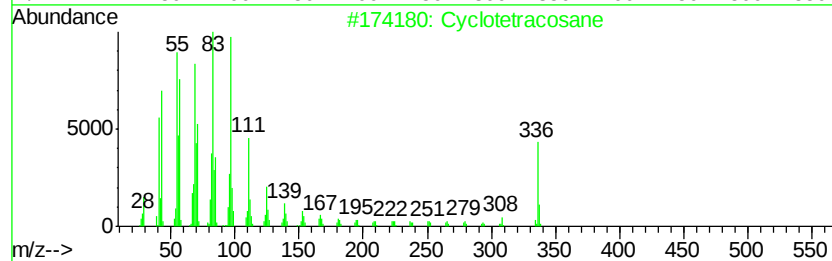
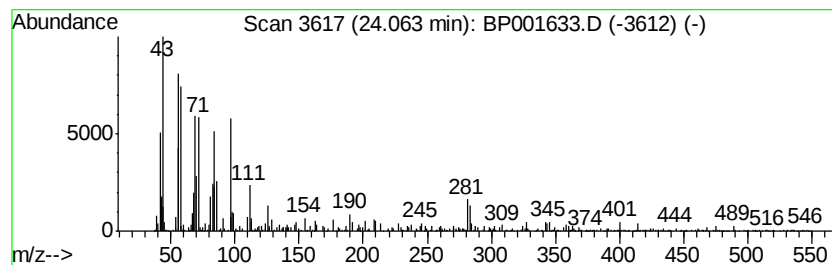
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Cyclotetracosane Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.06	2.65 ng	79170	Perylene-d12	23.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclotetracosane	336	C24H48	000297-03-0	95
2		Cetene	224	C16H32	000629-73-2	93
3		Cyclopentadecane	210	C15H30	000295-48-7	86
4		17-Pentatriacontene	491	C35H70	006971-40-0	83
5		1-Heneicosanol	312	C21H44O	015594-90-8	81



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP011620\
Data File : BP001633.D
Acq On : 16 Jan 2020 18:09
Operator : JU
Sample : L1145-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

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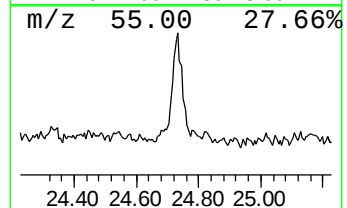
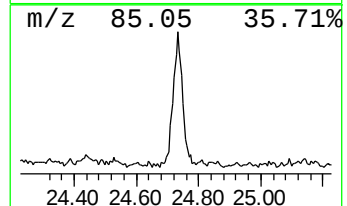
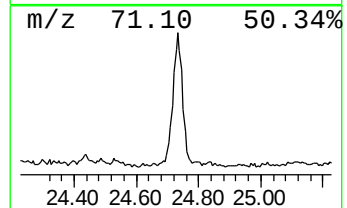
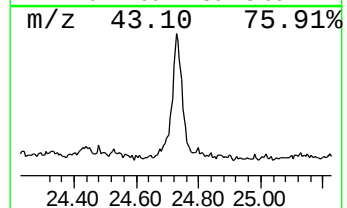
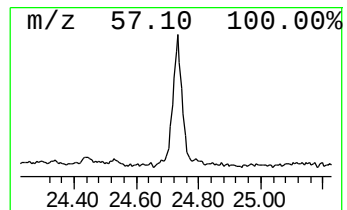
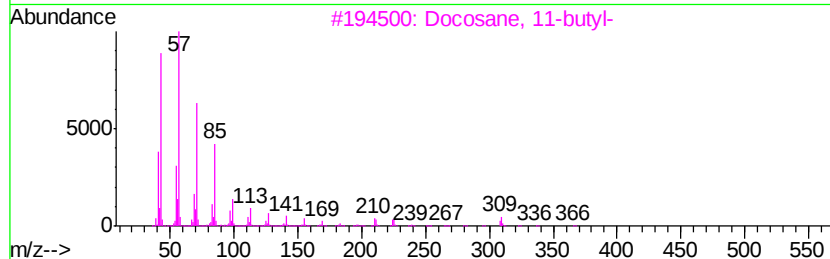
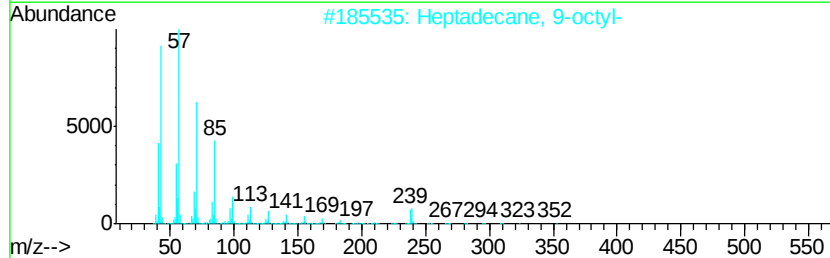
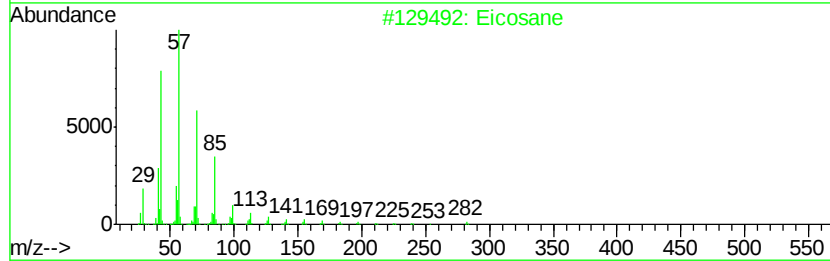
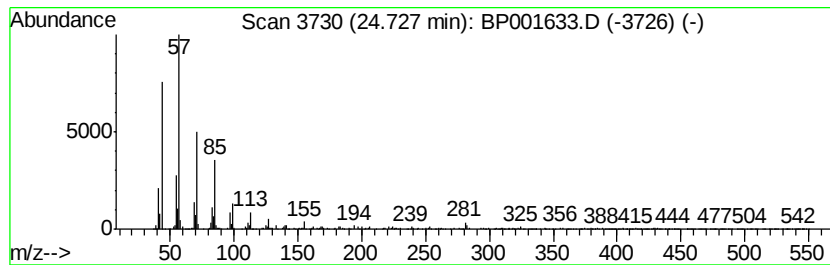
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 Docosane, 11-butyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.73	6.09 ng	182120	Perylene-d12	23.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	95
2		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	93
3		Docosane, 11-butyl-	366	C26H54	013475-76-8	91
4		Heneicosane, 11-pentyl-	366	C26H54	014739-72-1	90
5		Tetracosane	338	C24H50	000646-31-1	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP011620\
 Data File : BP001633.D
 Acq On : 16 Jan 2020 18:09
 Operator : JU
 Sample : L1145-01
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP010820.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
Butane, 2-methoxy...	2.95	49.4	ng	766387	1	7.63	310591	20.0
2-Pentanone, 4-hy...	4.79	5.1	ng	79561	1	7.63	310591	20.0
unknown7.17	7.17	74.8	ng	1160940	1	7.63	310591	20.0
2,4,7,9-Tetrameth...	13.20	2.3	ng	64294	3	14.27	553348	20.0
Tridecanoic acid	17.97	6.2	ng	204080	4	17.04	662713	20.0
1-Hexacosene	20.89	13.8	ng	441019	5	21.14	639735	20.0
Heneicosane, 11-d...	21.32	8.0	ng	256819	5	21.14	639735	20.0
Heneicosane	21.76	9.1	ng	290410	5	21.14	639735	20.0
Hexacosane	22.23	11.7	ng	373698	5	21.14	639735	20.0
Supraene	22.36	14.3	ng	427989	6	23.37	597684	20.0
2-Bromo dodecane	22.74	12.7	ng	379476	6	23.37	597684	20.0
2-methyloctacosane	23.32	13.6	ng	407410	6	23.37	597684	20.0
Eicosane, 7-hexyl-	23.97	11.9	ng	356006	6	23.37	597684	20.0
Cyclotetracosane	24.06	2.6	ng	79170	6	23.37	597684	20.0
Docosane, 11-butyl-	24.73	6.1	ng	182120	6	23.37	597684	20.0