

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP122019\
 Data File : BP001340.D
 Acq On : 20 Dec 2019 13:01
 Operator : JU
 Sample : K6300-01
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SSI-MW-4

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-BP121919.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	5.610	590	599	610	rBV	92066	134324	5.10%	0.733%
2	6.204	690	700	716	rBV	732520	919948	34.92%	5.020%
3	7.746	945	962	970	rBV	512684	640188	24.30%	3.494%
4	7.940	988	995	1006	rBV	1506307	1783704	67.70%	9.734%
5	8.298	1051	1056	1063	rBV	366656	411373	15.61%	2.245%
6	8.545	1091	1098	1102	rBV	1333057	1645623	62.46%	8.980%
7	9.187	1199	1207	1210	rBV	1114233	1397729	53.05%	7.628%
8	10.334	1396	1402	1410	rVB	420590	481280	18.27%	2.626%
9	12.116	1698	1705	1710	rBV	1992837	2423811	91.99%	13.227%
10	12.786	1814	1819	1825	rBV	62511	72016	2.73%	0.393%
11	13.198	1883	1889	1897	rVB	457269	549747	20.86%	3.000%
12	14.492	2102	2109	2121	rBV	1375529	1718380	65.22%	9.377%
13	15.598	2291	2297	2309	rBV	500781	564284	21.42%	3.079%
14	16.486	2444	2448	2463	rBV2	61030	91317	3.47%	0.498%
15	17.574	2630	2633	2641	rBV2	17208	27420	1.04%	0.150%
16	17.780	2664	2668	2672	rBV	23723	26884	1.02%	0.147%
17	17.886	2680	2686	2690	rBV	2387987	2634791	100.00%	14.378%
18	18.257	2745	2749	2754	rVB	66513	64750	2.46%	0.353%
19	18.704	2821	2825	2829	rVB	76739	77278	2.93%	0.422%
20	19.127	2891	2897	2911	rBV3	120656	224478	8.52%	1.225%
21	19.245	2911	2917	2922	rBV	406585	479392	18.19%	2.616%
22	19.539	2961	2967	2974	rVB	107611	122823	4.66%	0.670%
23	19.927	3028	3033	3041	rBV	172449	193692	7.35%	1.057%
24	20.304	3092	3097	3101	rBV	228676	260002	9.87%	1.419%
25	20.698	3159	3164	3171	rBV	219884	280598	10.65%	1.531%
26	21.015	3212	3218	3225	rVB	330479	447009	16.97%	2.439%
27	21.133	3232	3238	3245	rBV	180743	275382	10.45%	1.503%
28	21.633	3317	3323	3329	rBV	129168	216367	8.21%	1.181%
29	22.204	3414	3420	3427	rVB	60531	115416	4.38%	0.630%
30	22.874	3526	3534	3541	rBV2	18930	44556	1.69%	0.243%

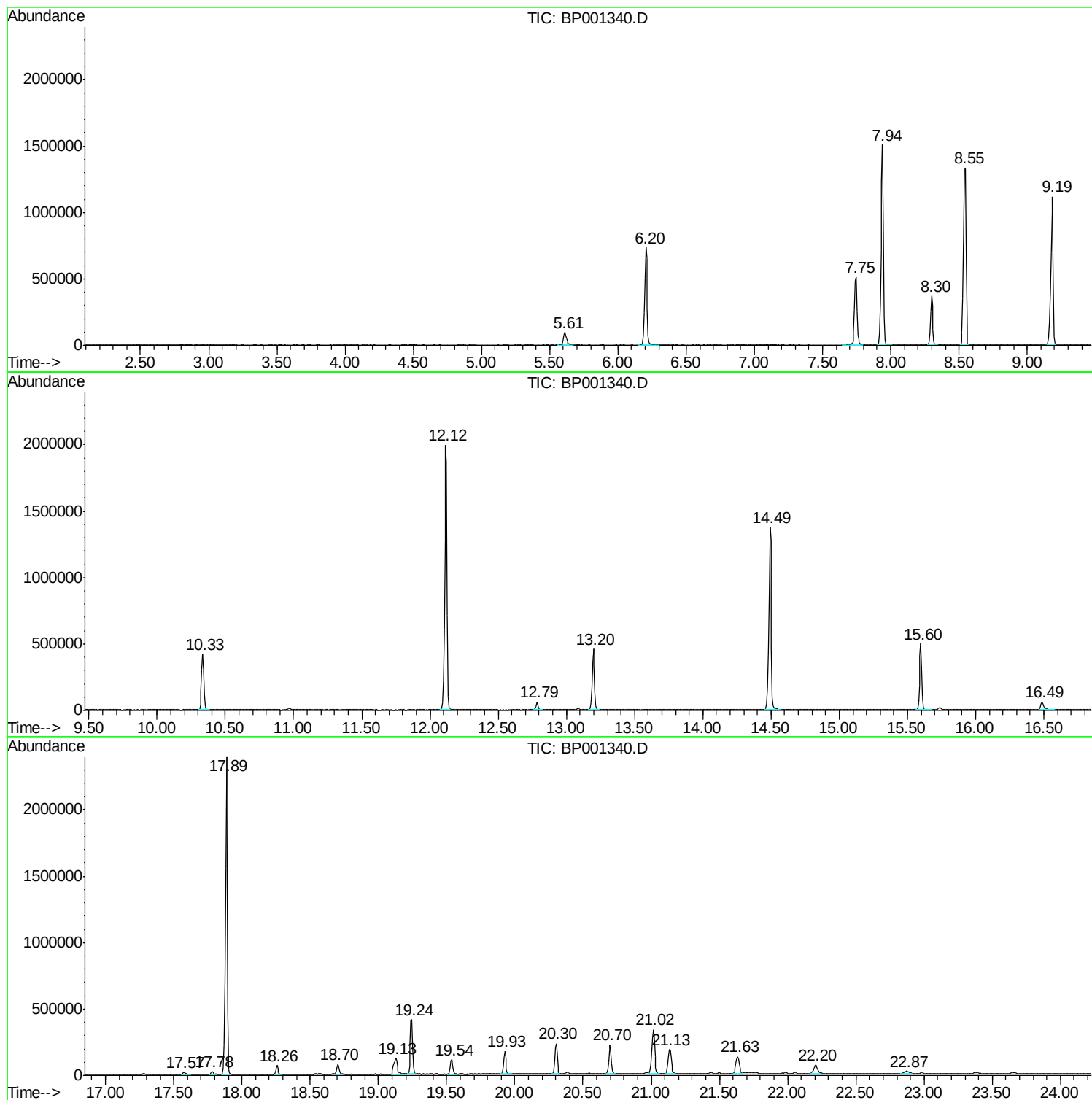
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP121919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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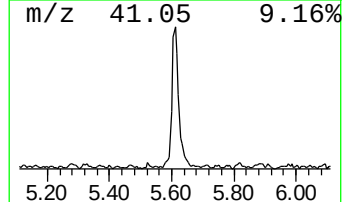
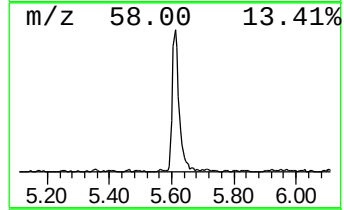
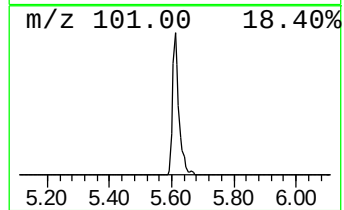
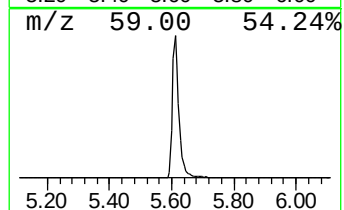
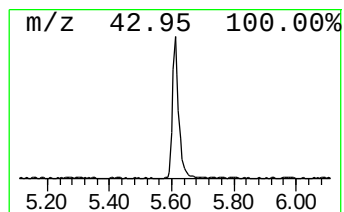
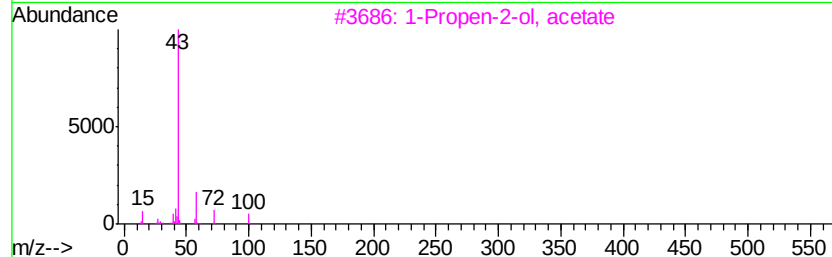
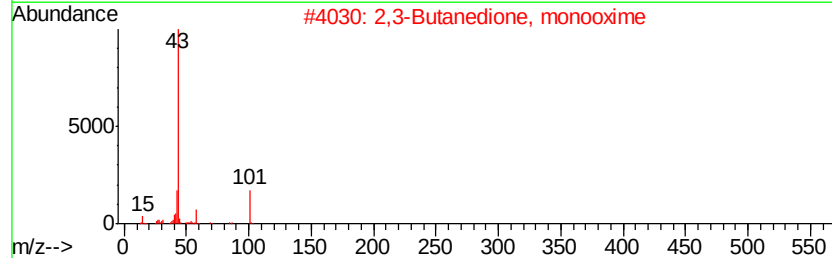
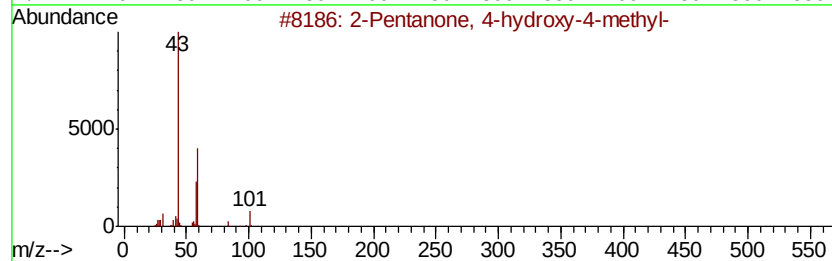
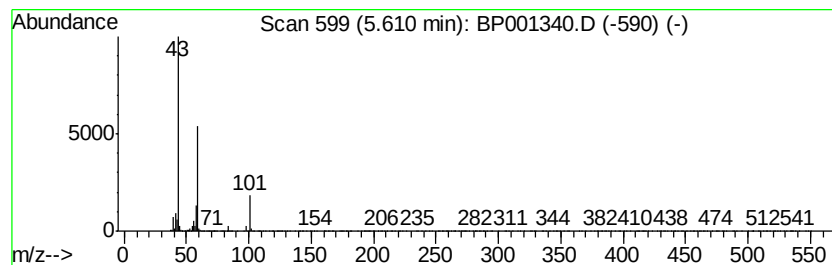
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.61	6.53 ng	134324	1,4-Dichlorobenzene-d4	8.30

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45
2	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16
3	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
4	Butane	58	C4H10	000106-97-8	9
5	Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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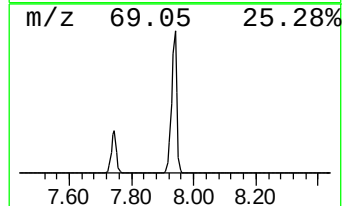
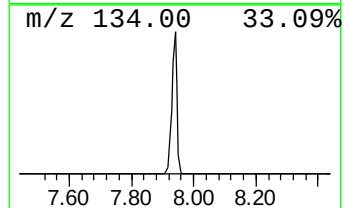
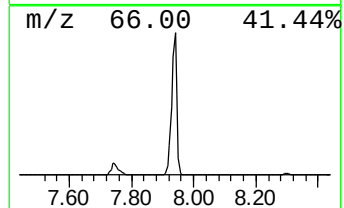
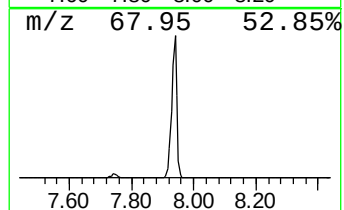
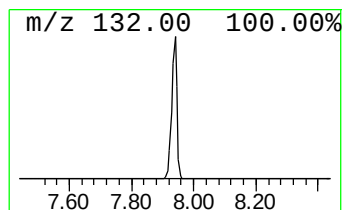
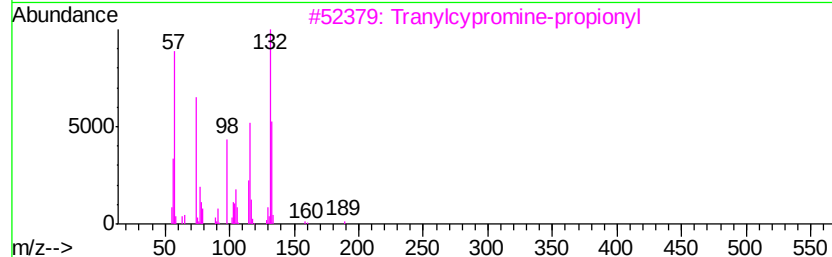
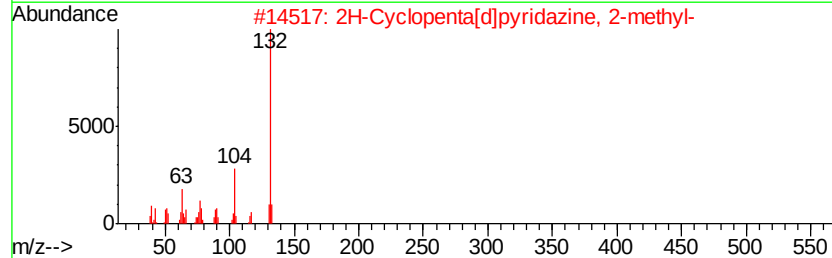
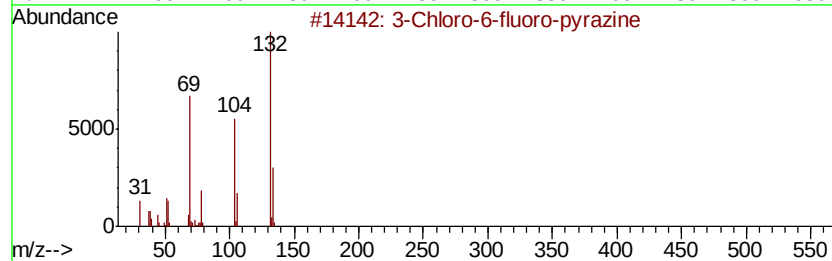
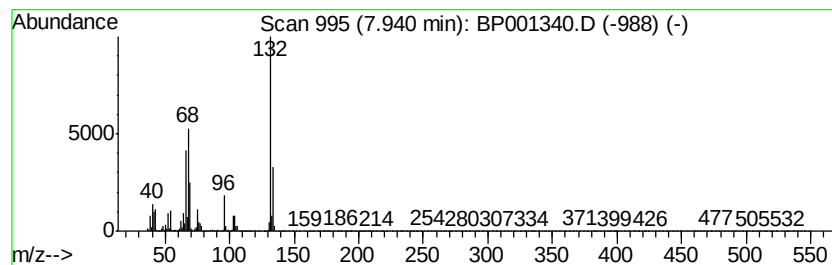
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown7.94 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.94	86.72 ng	1783700	1,4-Dichlorobenzene-d4	8.30

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	14
2		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	10
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	10
4		Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	10
5		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9



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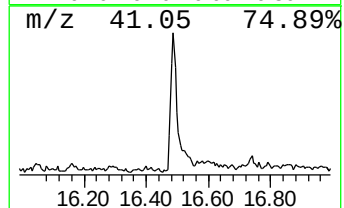
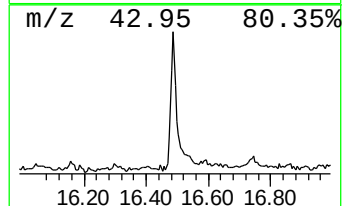
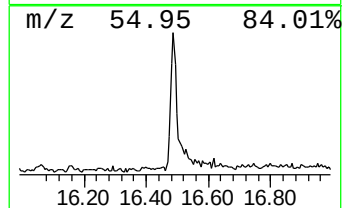
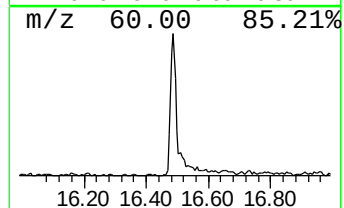
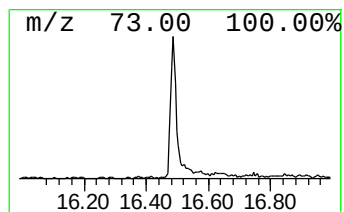
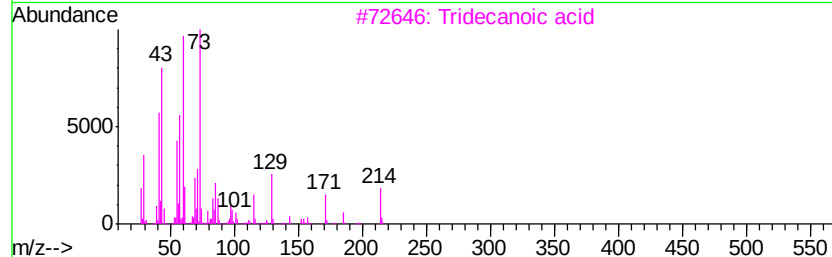
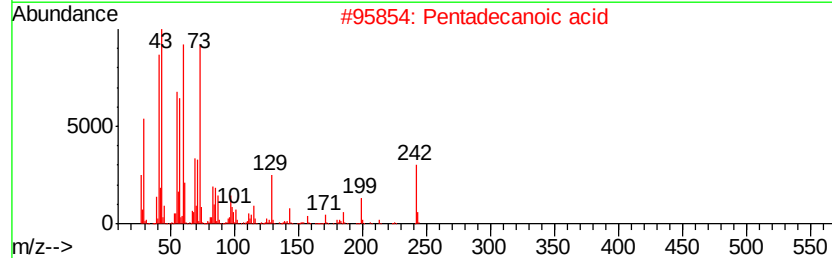
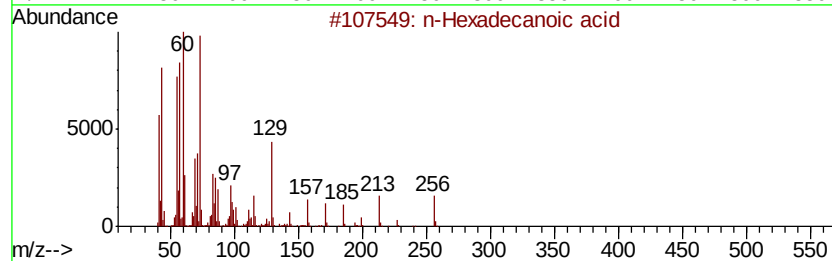
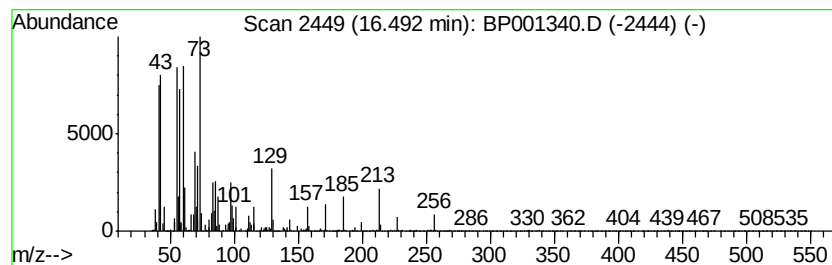
Instrument :
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 ClientSampleId :
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TIC Library : C:\DATABASE\NIST11.L
 TIC Integration Parameters: LSCINT.P

 Peak Number 4 n-Hexadecanoic acid Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD		R.T.	
16.49	3.24 ng	91317	Phenanthrene-d10		15.60	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
2		Pentadecanoic acid	242	C15H30O2	001002-84-2	87
3		Tridecanoic acid	214	C13H26O2	000638-53-9	76
4		Tetradecanoic acid	228	C14H28O2	000544-63-8	76
5		Octadecanoic acid	284	C18H36O2	000057-11-4	74



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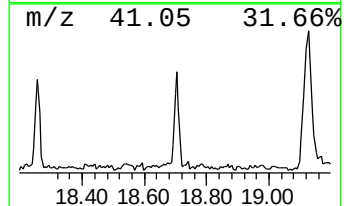
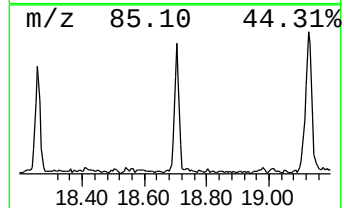
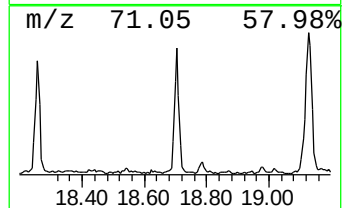
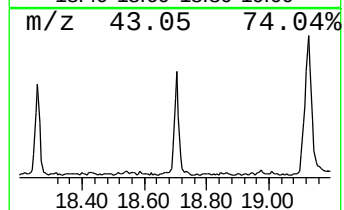
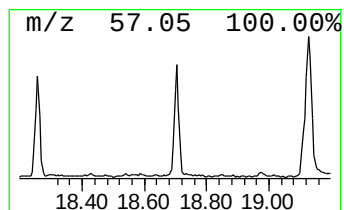
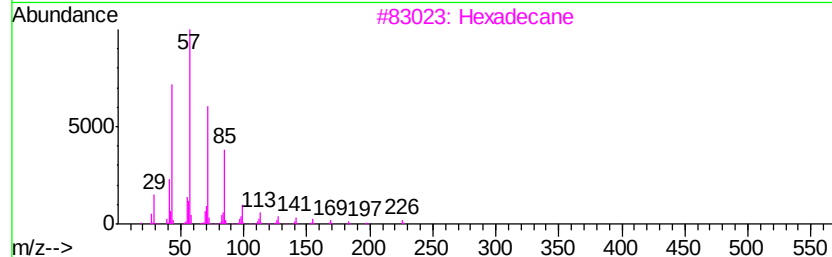
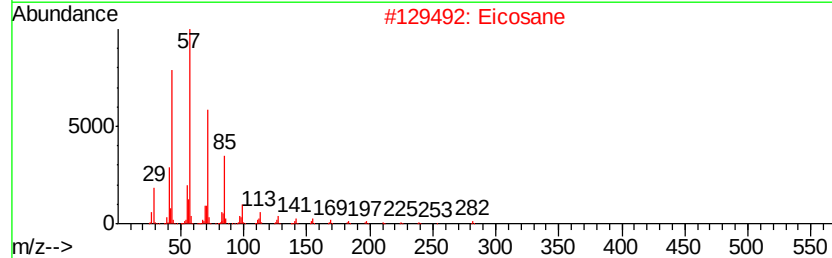
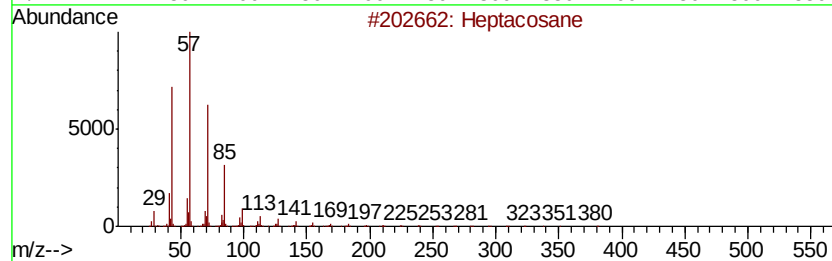
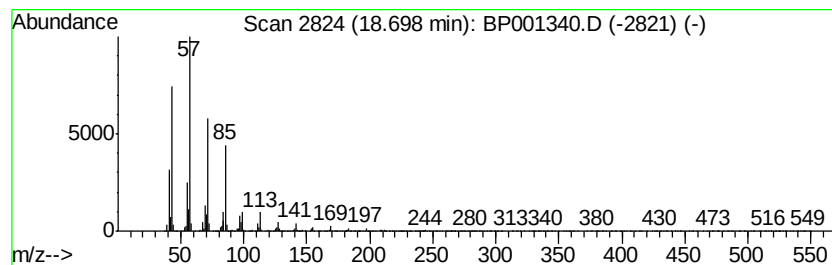
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Heptacosane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.70	3.22 ng	77278	Chrysene-d12	19.24
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Heptacosane	380 C27H56	000593-49-7	91
2	Eicosane	282 C20H42	000112-95-8	91
3	Hexadecane	226 C16H34	000544-76-3	91
4	Heneicosane	296 C21H44	000629-94-7	90
5	Pentadecane	212 C15H32	000629-62-9	90



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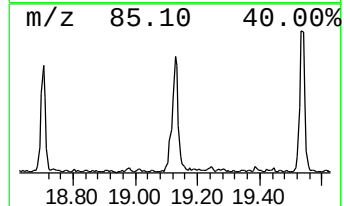
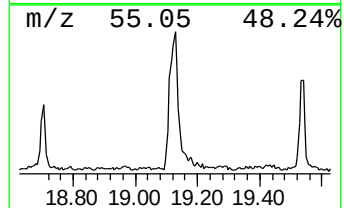
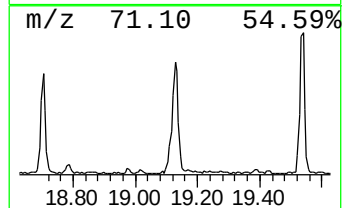
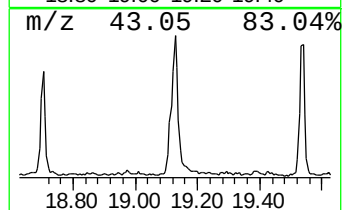
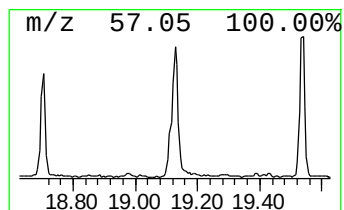
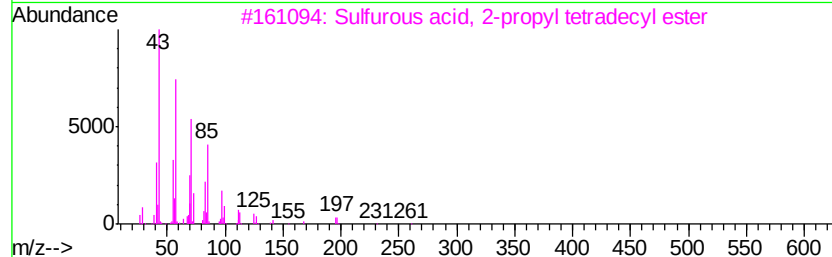
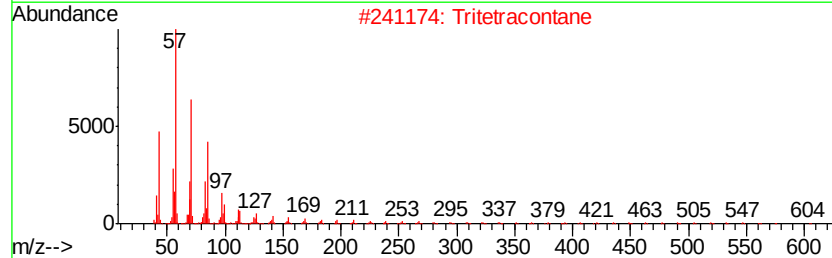
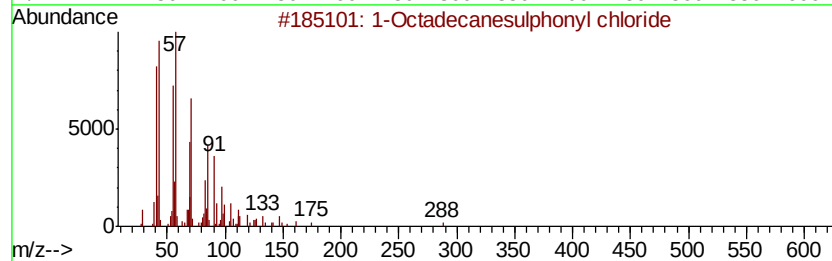
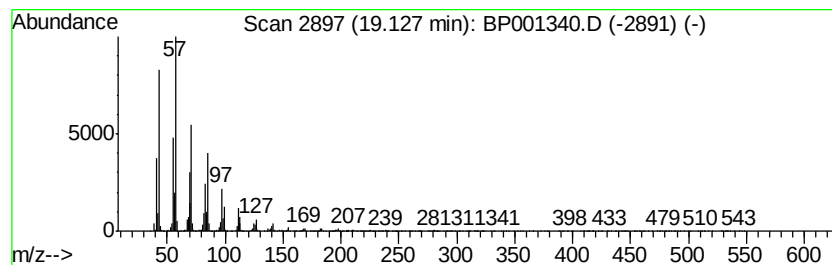
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 1-Octadecanesulphonyl chloride Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.13	9.37 ng	224478	Chrysene-d12	19.24

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Octadecanesulphonyl chloride	352	C18H37ClO2S	1000342-70-4	91
2		Tritetracontane	605	C43H88	007098-21-7	91
3		Sulfurous acid, 2-propyl tetradecyl	320	C17H36O3S	1000309-12-5	87
4		Sulfurous acid, octadecyl 2-propyl	376	C21H44O3S	1000309-12-7	83
5		Tetracosane, 1-bromo-	416	C24H49Br	006946-24-3	81



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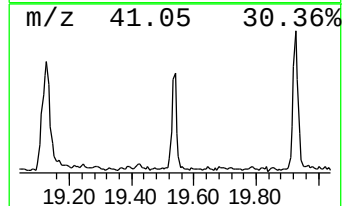
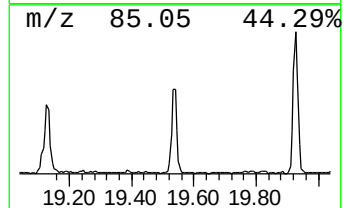
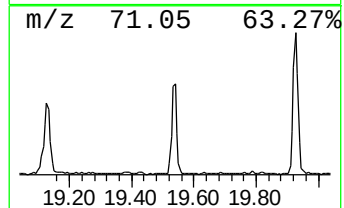
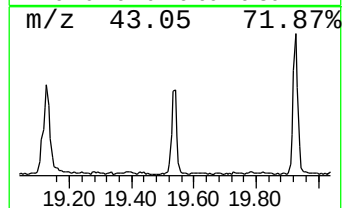
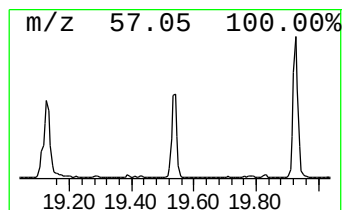
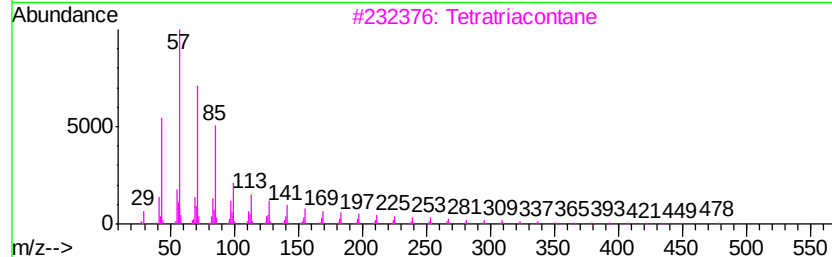
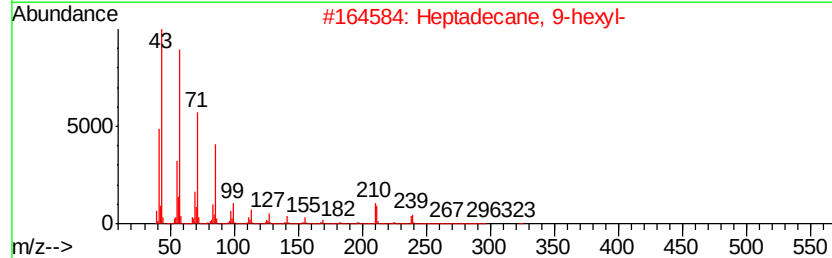
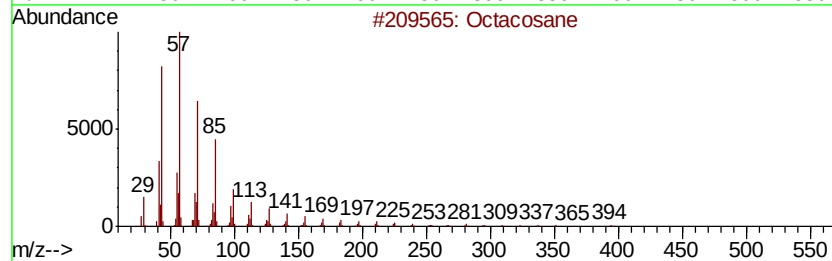
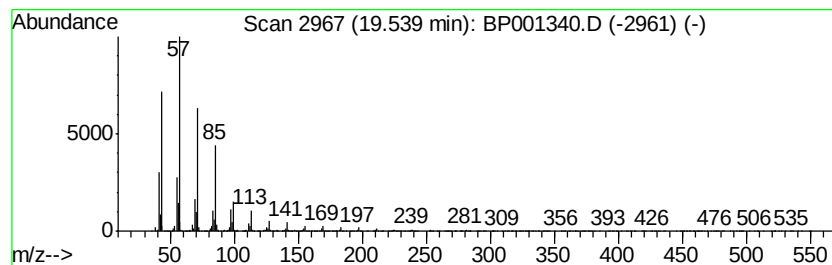
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP121919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Octacosane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.54	5.12 ng	122823	Chrysene-d12	19.24
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Octacosane		394 C28H58	000630-02-4 98
2	Heptadecane, 9-hexyl-		324 C23H48	055124-79-3 94
3	Tetratriacontane		479 C34H70	014167-59-0 93
4	Eicosane, 10-methyl-		296 C21H44	054833-23-7 93
5	Heptacosane		380 C27H56	000593-49-7 91



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP122019\
Data File : BP001340.D
Acq On : 20 Dec 2019 13:01
Operator : JU
Sample : K6300-01
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SSI-MW-4

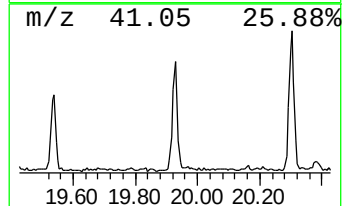
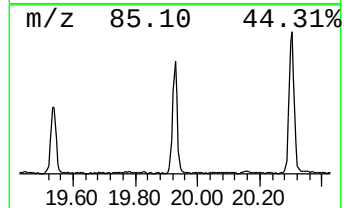
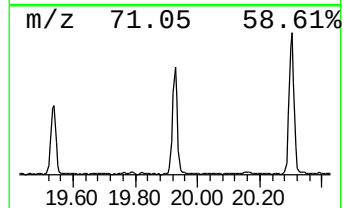
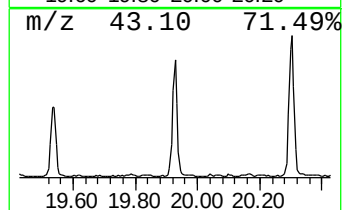
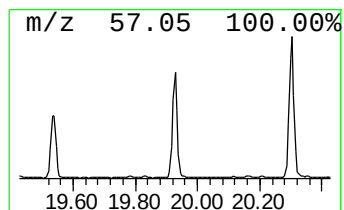
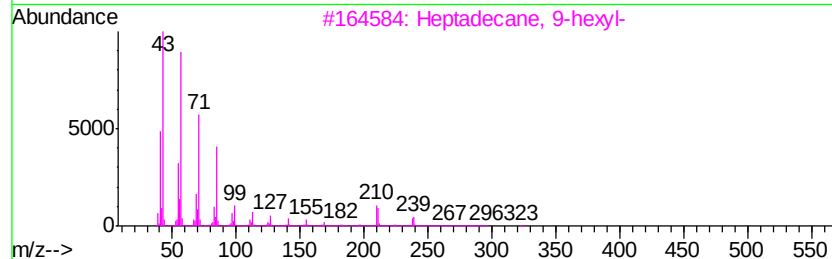
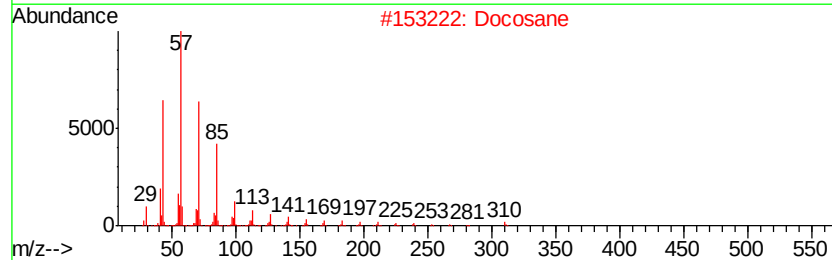
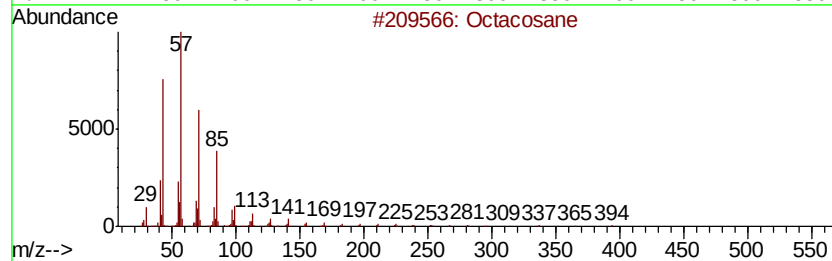
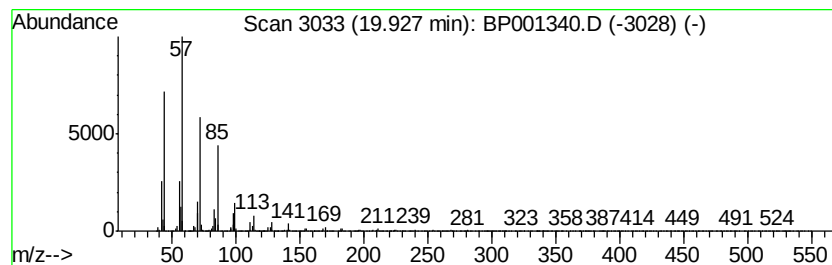
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Docosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.93	8.08 ng	193692	Chrysene-d12	19.24

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octacosane		394	C28H58	000630-02-4	95
2	Docosane		310	C22H46	000629-97-0	93
3	Heptadecane, 9-hexyl-		324	C23H48	055124-79-3	93
4	2-Bromo dodecane		248	C12H25Br	013187-99-0	91
5	Hexacosane		366	C26H54	000630-01-3	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP122019\
Data File : BP001340.D
Acq On : 20 Dec 2019 13:01
Operator : JU
Sample : K6300-01
Misc :
ALS Vial : 7 Sample Multiplier: 1

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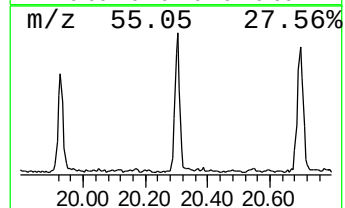
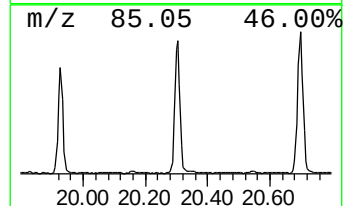
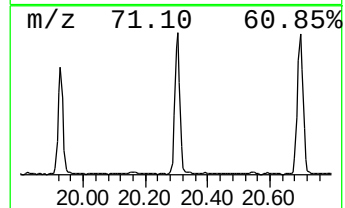
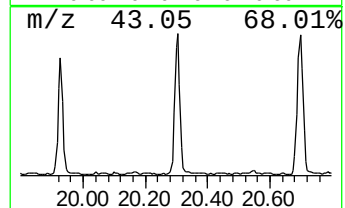
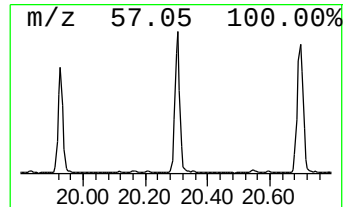
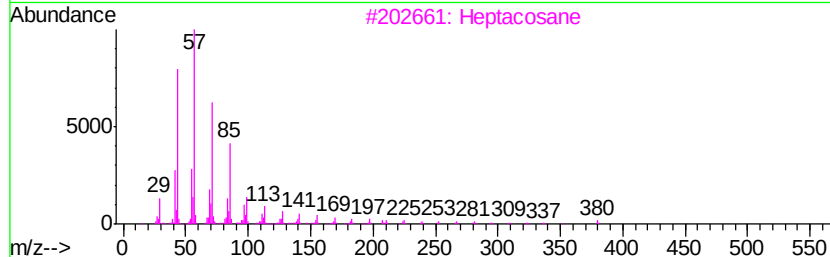
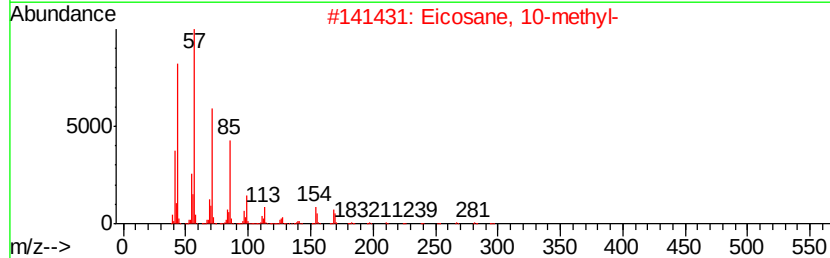
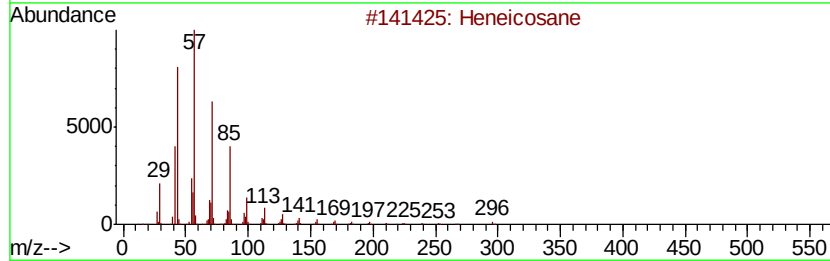
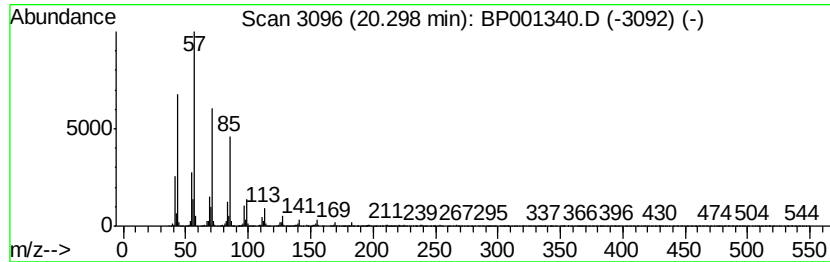
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Eicosane, 10-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.30	11.63 ng	260002	Perylene-d12	21.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	96
2		Eicosane, 10-methyl-	296	C21H44	054833-23-7	95
3		Heptacosane	380	C27H56	000593-49-7	95
4		Eicosane, 3-methyl-	296	C21H44	006418-46-8	94
5		Hexacosane	366	C26H54	000630-01-3	94



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP122019\
Data File : BP001340.D
Acq On : 20 Dec 2019 13:01
Operator : JU
Sample : K6300-01
Misc :
ALS Vial : 7 Sample Multiplier: 1

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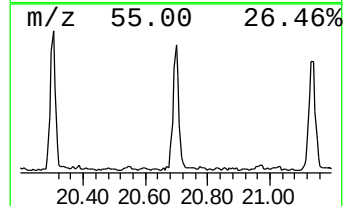
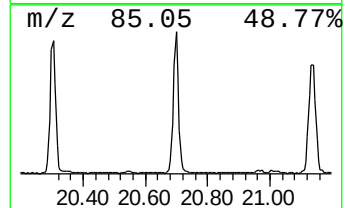
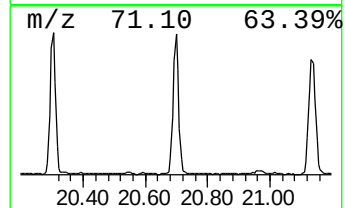
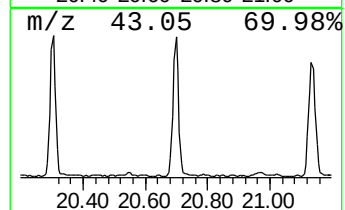
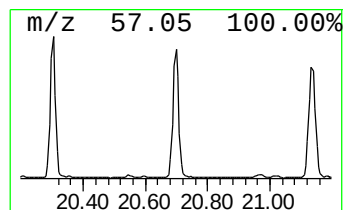
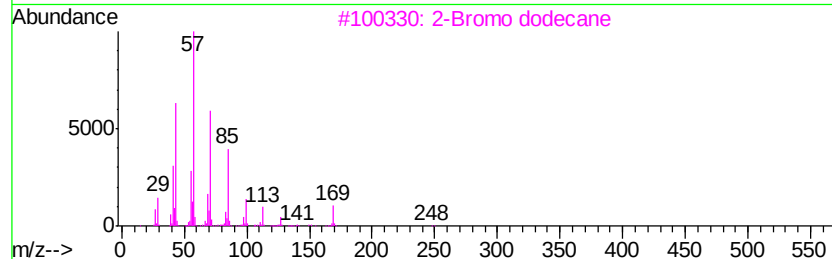
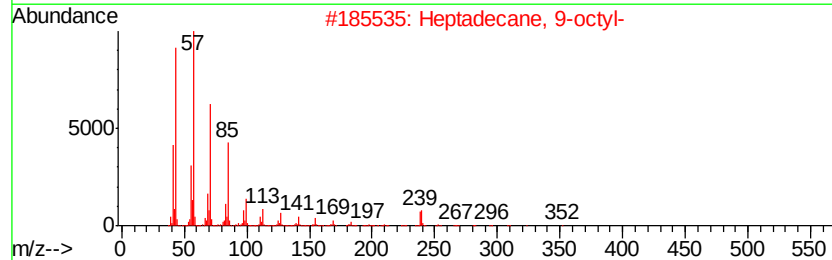
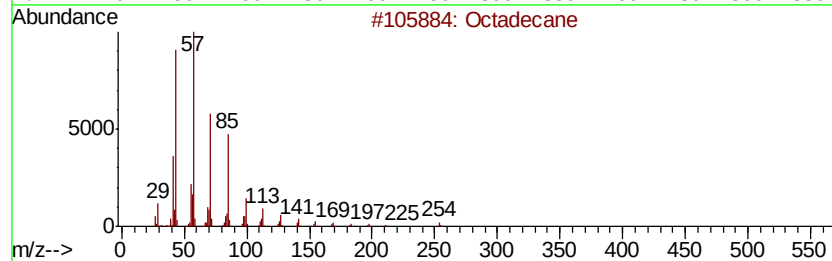
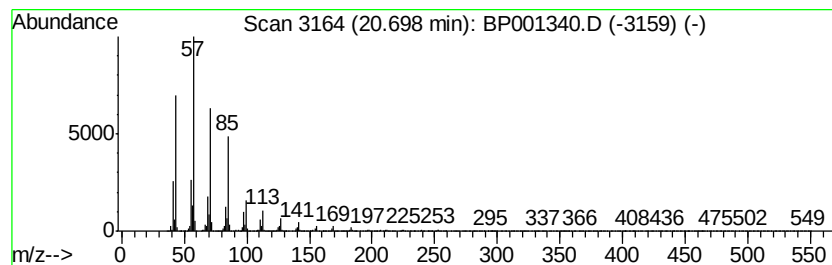
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Heptadecane, 9-octyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.70	12.55 ng	280598	Perylene-d12	21.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecane	254	C18H38	000593-45-3	94
2		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	93
3		2-Bromo dodecane	248	C12H25Br	013187-99-0	91
4		Eicosane, 3-methyl-	296	C21H44	006418-46-8	91
5		Octacosane	394	C28H58	000630-02-4	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP122019\
Data File : BP001340.D
Acq On : 20 Dec 2019 13:01
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Misc :
ALS Vial : 7 Sample Multiplier: 1

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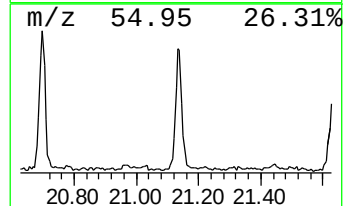
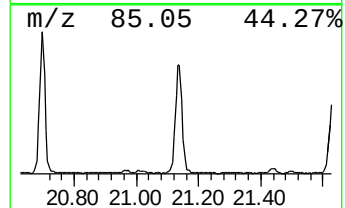
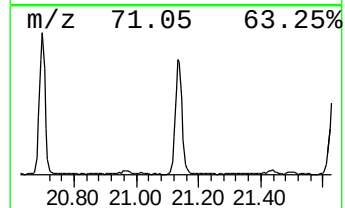
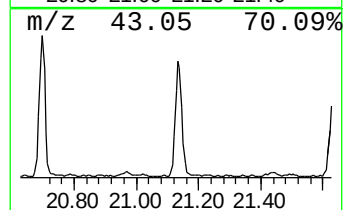
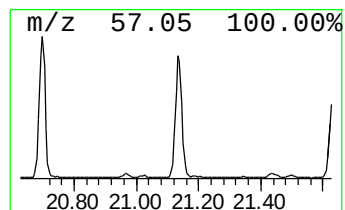
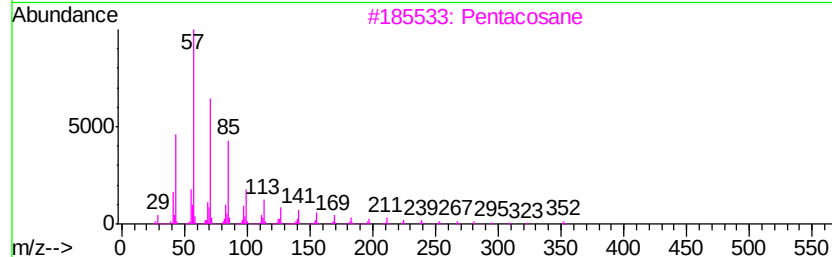
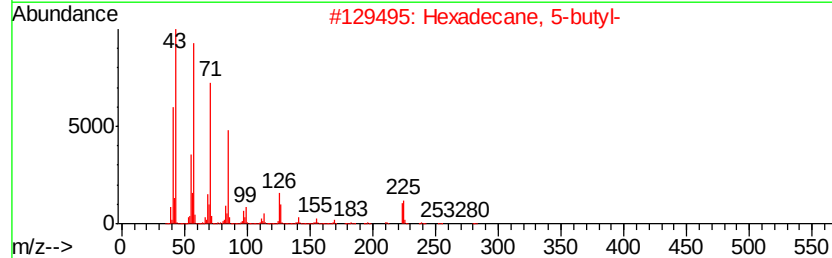
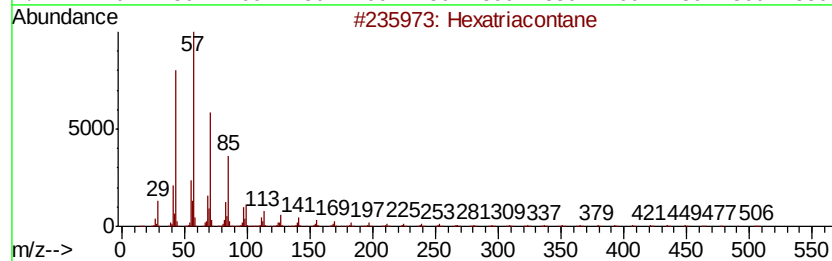
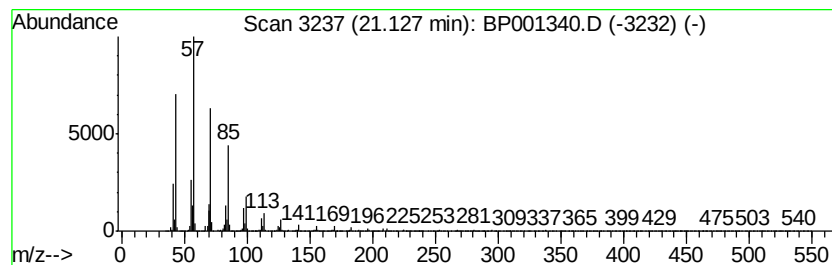
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Hexatriacontane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.13	12.32 ng	275382	Perylene-d12	21.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexatriacontane	507	C36H74	000630-06-8	91
2		Hexadecane, 5-butyl-	282	C20H42	006912-07-8	91
3		Pentacosane	352	C25H52	000629-99-2	91
4		Heneicosane	296	C21H44	000629-94-7	91
5		Hexacosane	366	C26H54	000630-01-3	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP122019\
 Data File : BP001340.D
 Acq On : 20 Dec 2019 13:01
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 Misc :
 ALS Vial : 7 Sample Multiplier: 1

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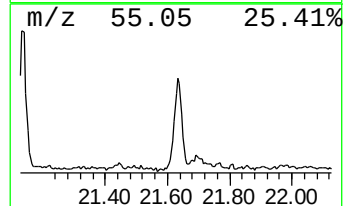
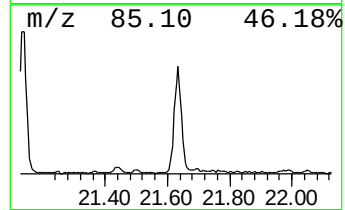
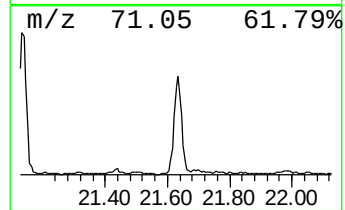
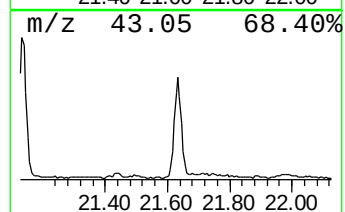
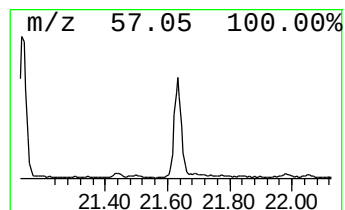
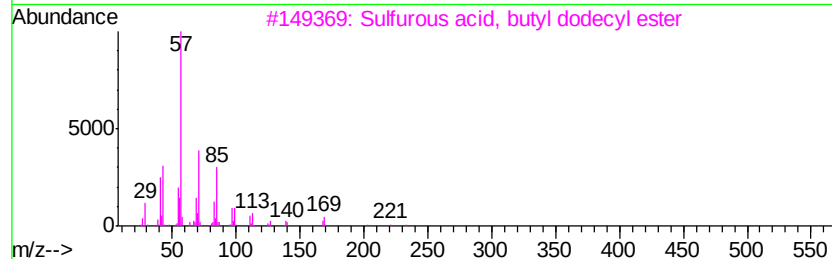
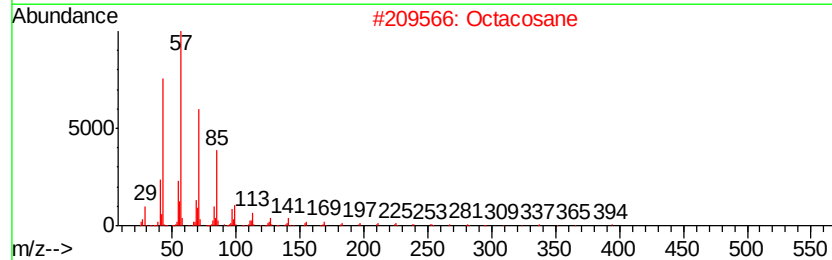
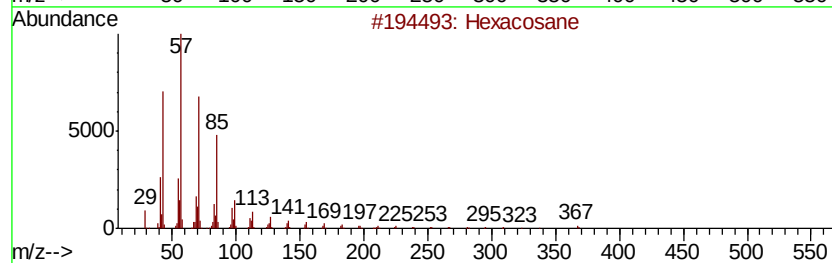
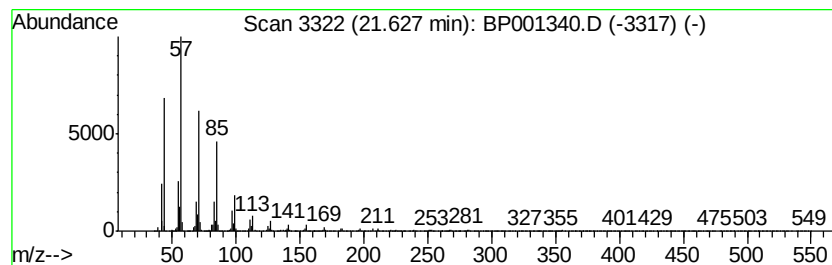
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 13 Sulfurous acid, butyl dodec... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.63	9.68 ng	216367	Perylene-d12	21.02

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexacosane	366	C26H54	000630-01-3	91
2			Octacosane	394	C28H58	000630-02-4	91
3			Sulfurous acid, butyl dodecyl ester	306	C16H34O3S	1000309-17-9	90
4			Tetratetracontane	619	C44H90	007098-22-8	87
5			Octadecane	254	C18H38	000593-45-3	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP122019\
Data File : BP001340.D
Acq On : 20 Dec 2019 13:01
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Misc :
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ClientSampleId :
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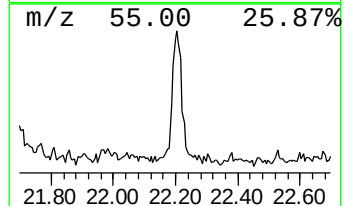
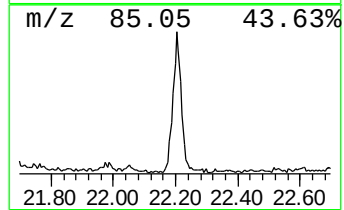
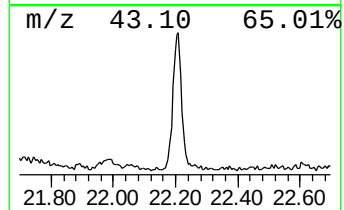
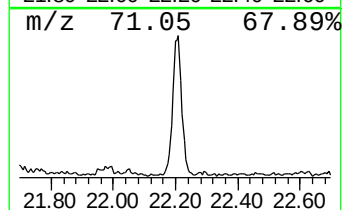
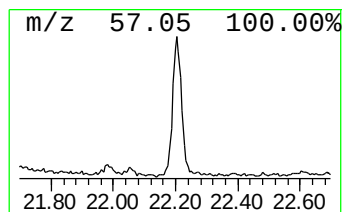
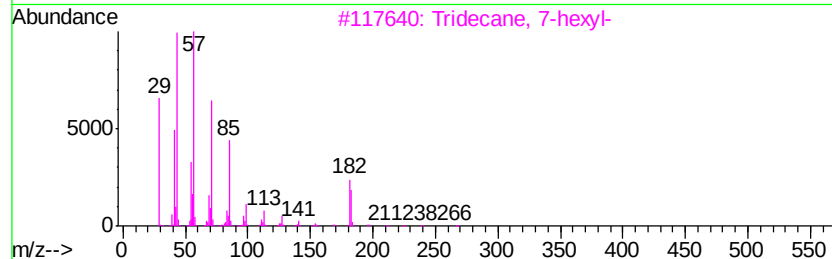
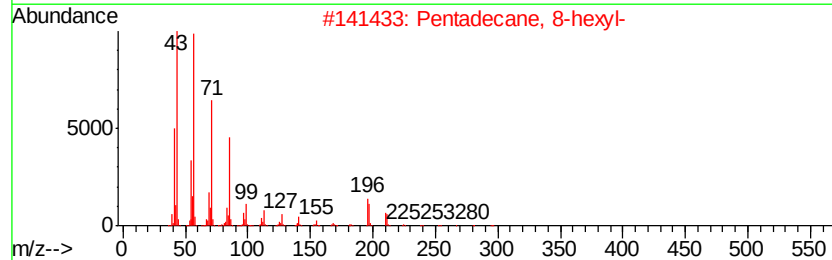
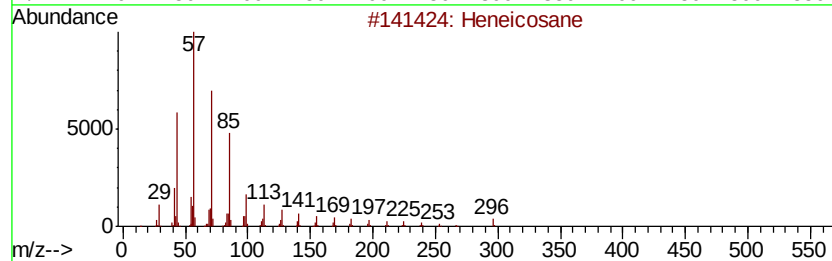
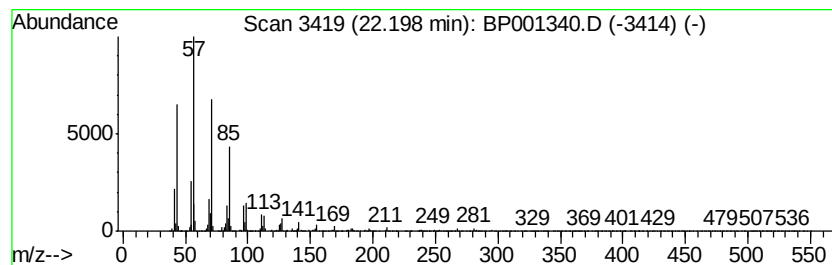
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Pentadecane, 8-hexyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.20	5.16 ng	115416	Perylene-d12	21.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	97
2		Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	94
3		Tridecane, 7-hexyl-	268	C19H40	007225-66-3	93
4		Tricosane	324	C23H48	000638-67-5	93
5		Octadecane, 5,14-dibutyl-	366	C26H54	055282-13-8	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP122019\
 Data File : BP001340.D
 Acq On : 20 Dec 2019 13:01
 Operator : JU
 Sample : K6300-01
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 ALS Vial : 7 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP121919.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
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unknown7.94	7.94	86.7	ng	1783700	1	8.30	411373	20.0
n-Hexadecanoic acid	16.49	3.2	ng	91317	4	15.60	564284	20.0
Heptacosane	18.70	3.2	ng	77278	5	19.24	479392	20.0
1-Octadecanesulph...	19.13	9.4	ng	224478	5	19.24	479392	20.0
Octacosane	19.54	5.1	ng	122823	5	19.24	479392	20.0
Docosane	19.93	8.1	ng	193692	5	19.24	479392	20.0
Eicosane, 10-methyl-	20.30	11.6	ng	260002	6	21.02	447009	20.0
Heptadecane, 9-oc...	20.70	12.6	ng	280598	6	21.02	447009	20.0
Hexatriacontane	21.13	12.3	ng	275382	6	21.02	447009	20.0
Sulfurous acid, b...	21.63	9.7	ng	216367	6	21.02	447009	20.0
Pentadecane, 8-he...	22.20	5.2	ng	115416	6	21.02	447009	20.0