

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010320\
 Data File : BG043904.D
 Acq On : 4 Jan 2020 1:48
 Operator : JU
 Sample : K6449-05
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 OW-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_G\METHODS\8270-BG123019.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	3.147	5	10	28	rVB	2264156	3369443	34.96%	5.793%
2	5.074	331	338	349	rVB	328429	517707	5.37%	0.890%
3	5.544	411	418	427	rBV	1632002	2613983	27.12%	4.494%
4	7.130	679	688	698	rBV	1227804	2137968	22.18%	3.676%
5	7.501	743	751	759	rBV	2564608	4411093	45.76%	7.584%
6	7.965	823	830	838	rBV	471600	791355	8.21%	1.361%
7	8.288	876	885	893	rBV	1852565	3268091	33.91%	5.619%
8	9.116	1018	1026	1037	rBV	1590827	2891344	30.00%	4.971%
9	9.580	1099	1105	1113	rBV	111925	196909	2.04%	0.339%
10	10.761	1298	1306	1315	rBV	659024	1204475	12.50%	2.071%
11	13.217	1717	1724	1739	rVB	4125279	6738266	69.91%	11.585%
12	14.046	1855	1865	1870	rBV	288619	448405	4.65%	0.771%
13	14.251	1895	1900	1906	rVB2	92840	145322	1.51%	0.250%
14	14.592	1952	1958	1966	rVV	1149969	1797336	18.65%	3.090%
15	16.079	2205	2211	2228	rBV	3654149	5831784	60.50%	10.026%
16	17.336	2418	2425	2434	rBV	1395522	2144244	22.25%	3.687%
17	18.247	2575	2580	2587	rBV2	125333	228995	2.38%	0.394%
18	19.956	2864	2871	2876	rBV	6897901	9638848	100.00%	16.572%
19	20.767	3006	3009	3013	rVB	122154	134929	1.40%	0.232%
20	21.255	3087	3092	3105	rBV	1154686	1800501	18.68%	3.096%
21	21.584	3143	3148	3156	rVB2	1306827	2138686	22.19%	3.677%
22	21.807	3182	3186	3193	rVB	260962	370332	3.84%	0.637%
23	22.407	3283	3288	3295	rBV2	321036	575660	5.97%	0.990%
24	23.088	3398	3404	3412	rVB	338826	638851	6.63%	1.098%
25	23.875	3532	3538	3546	rBV2	339496	693313	7.19%	1.192%
26	24.716	3672	3681	3690	rBV2	768139	2178166	22.60%	3.745%
27	24.810	3690	3697	3707	rVB2	282021	695048	7.21%	1.195%
28	25.914	3877	3885	3895	rVB2	212730	563545	5.85%	0.969%

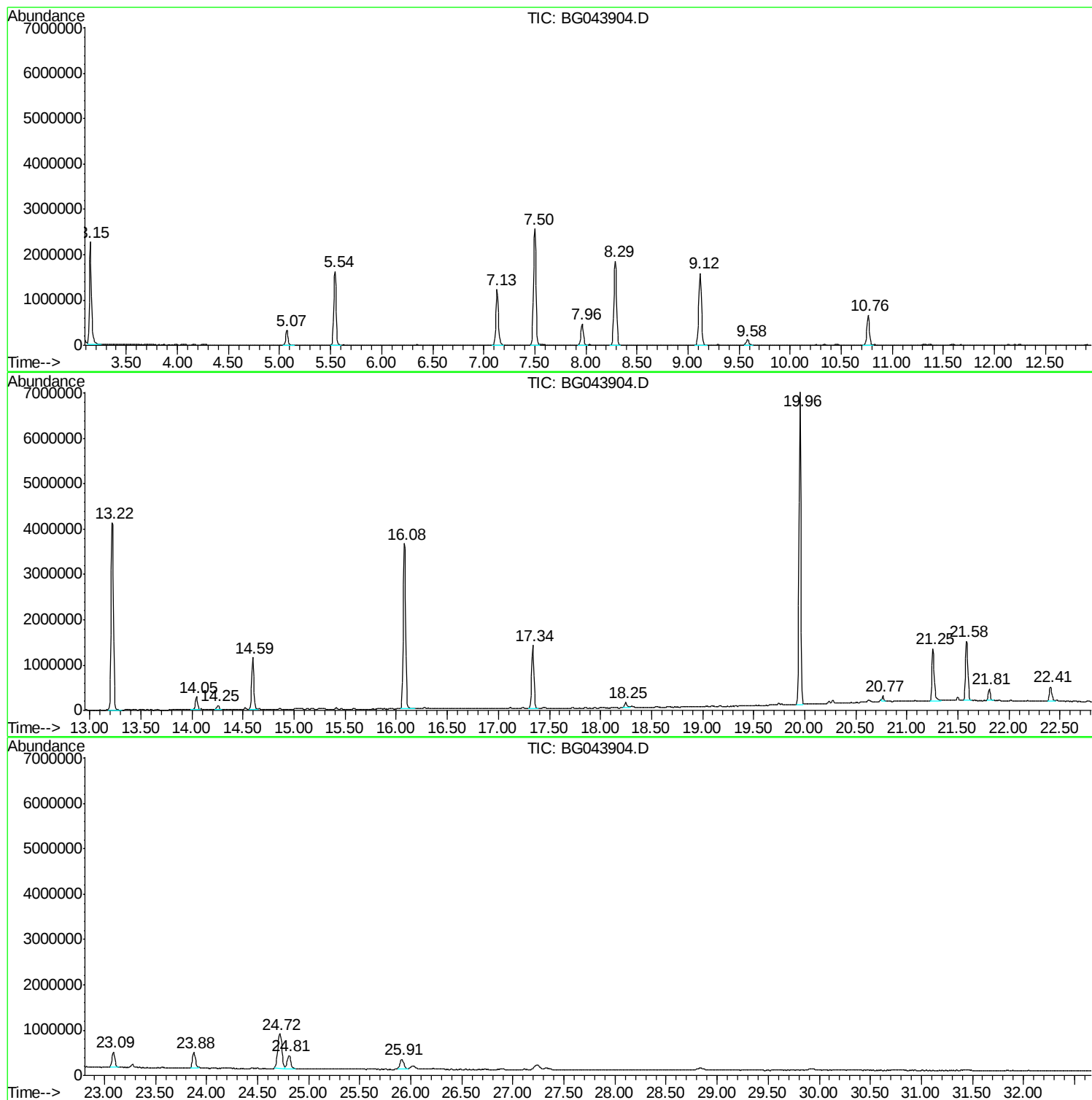
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG123019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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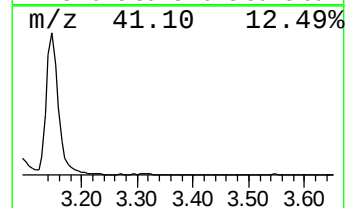
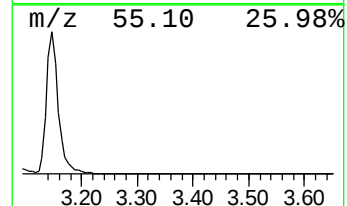
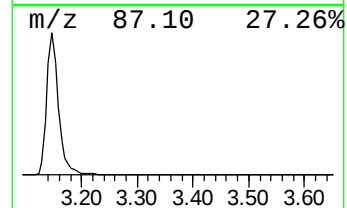
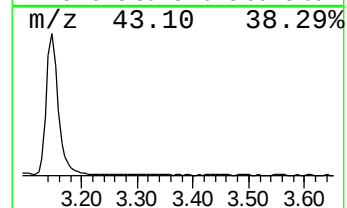
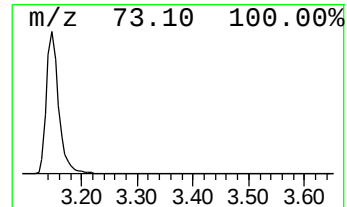
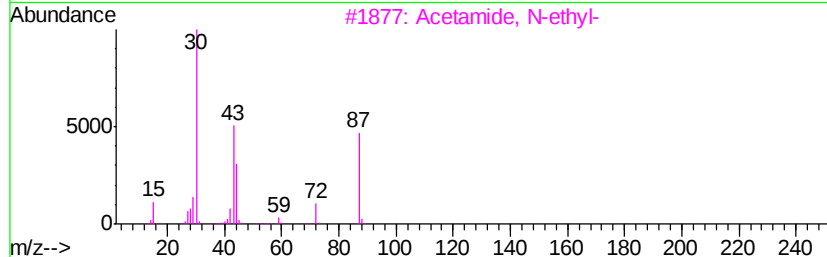
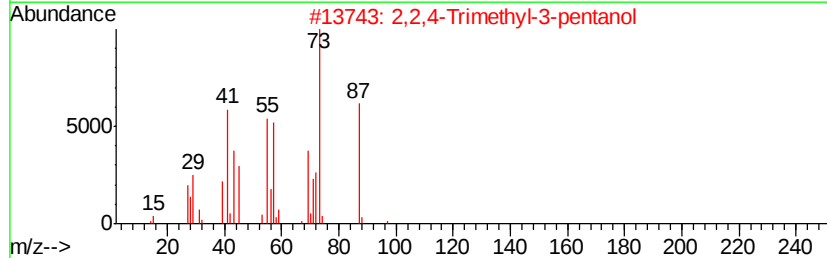
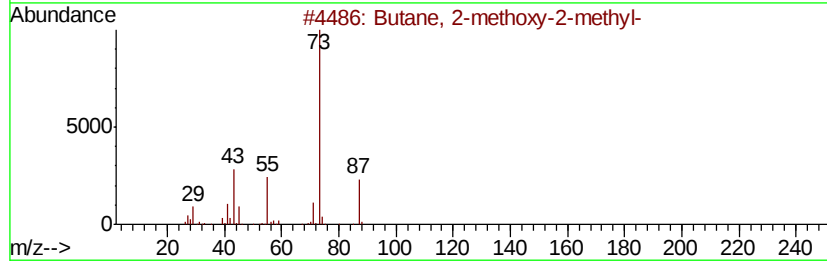
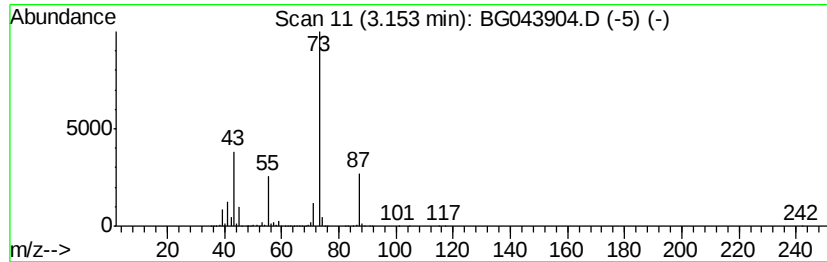
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.15	85.16 ng	3369440	1,4-Dichlorobenzene-d4	7.96

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	50
3		Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	35
4		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	25
5		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12



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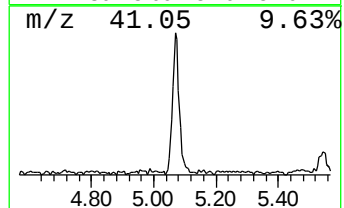
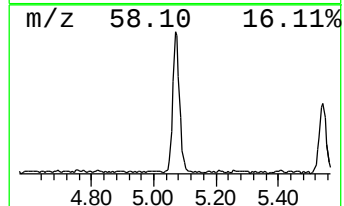
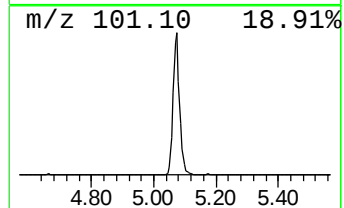
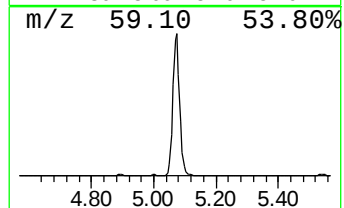
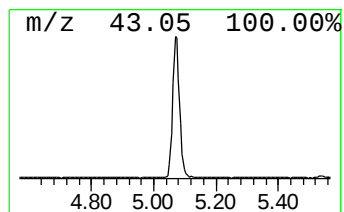
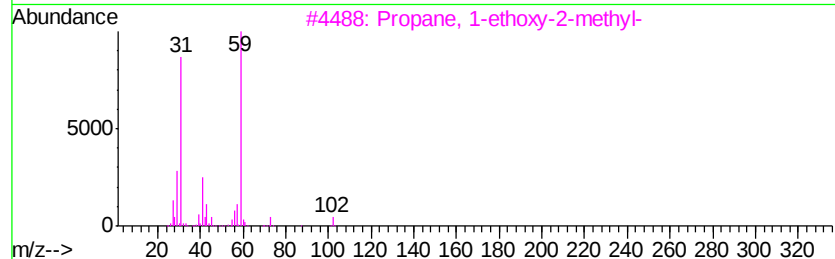
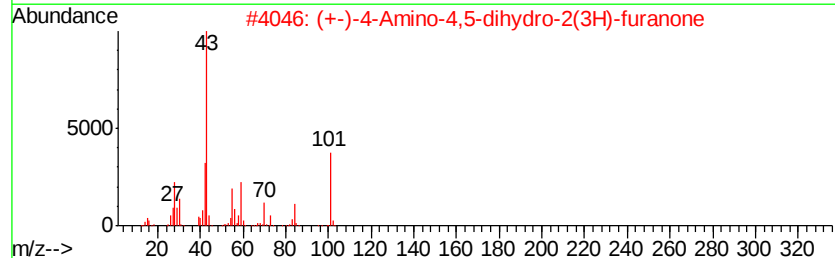
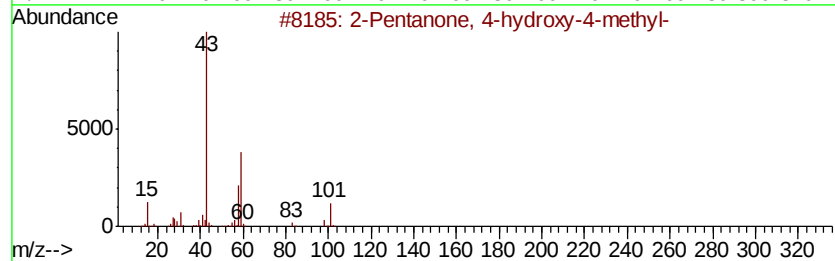
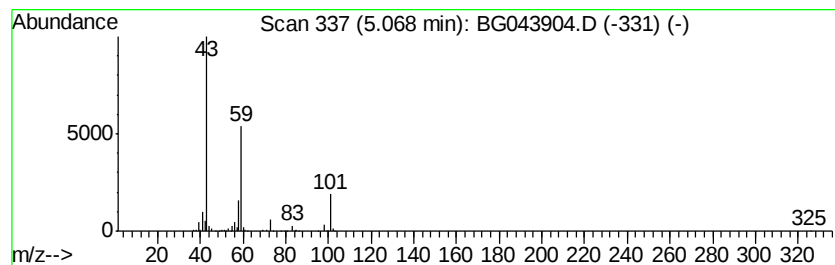
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.07	13.08 ng	517707	1,4-Dichlorobenzene-d4	7.96

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2		(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	23
3		Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	23
4		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	17
5		Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9



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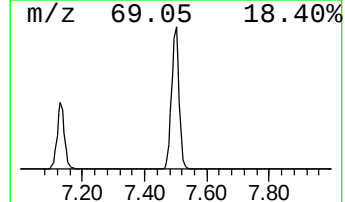
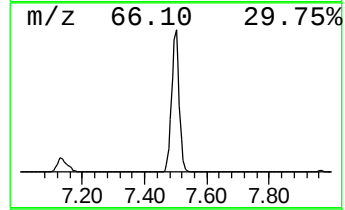
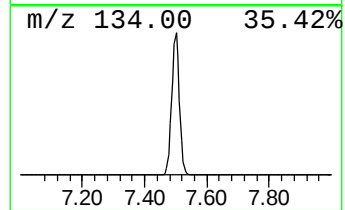
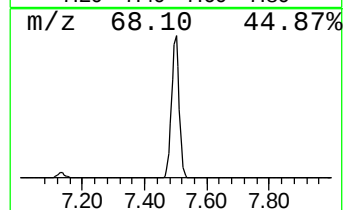
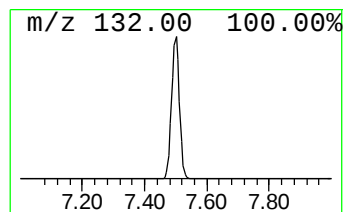
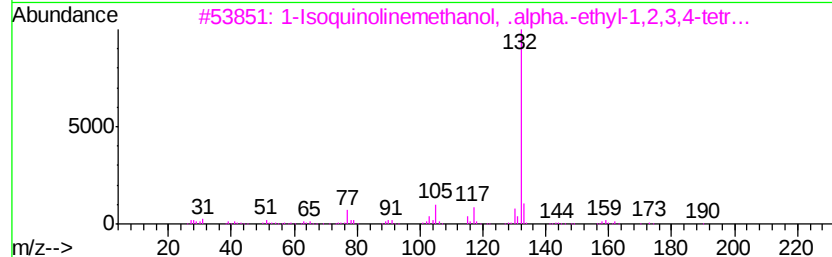
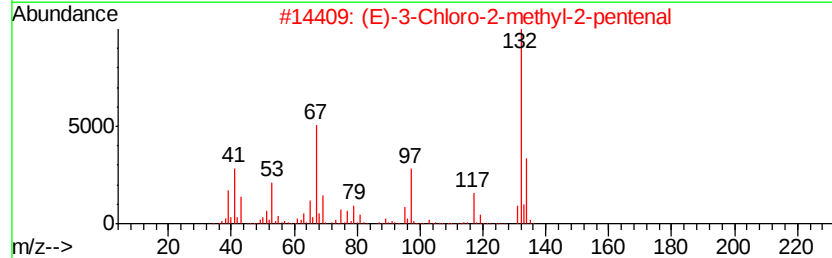
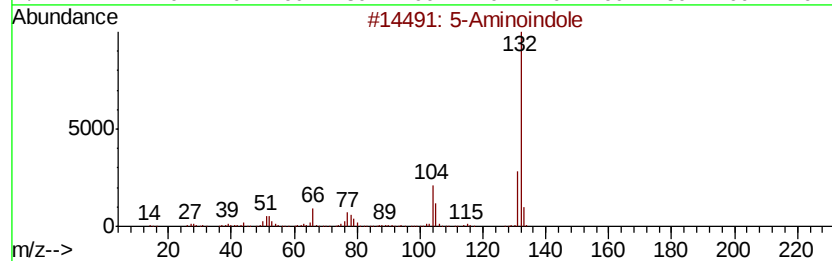
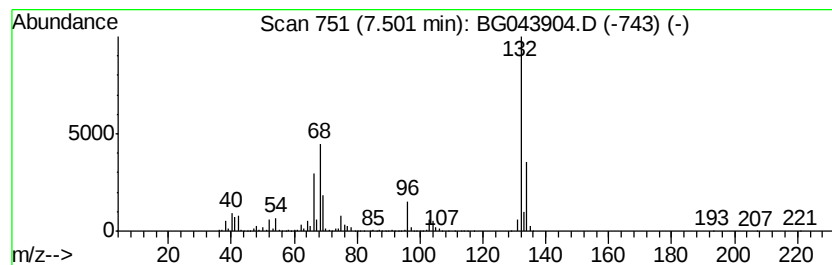
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Peak Number 3 unknown-01 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.50	111.48 ng	4411090	1,4-Dichlorobenzene-d4	7.96

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Aminoindole		132	C8H8N2	005192-03-0	22
2	(E)-3-Chloro-2-methyl-2-pentenal		132	C6H9ClO	031357-76-3	16
3	1-Isoquinolinemethanol, .alpha.-...		191	C12H17NO	114608-92-3	12
4	Tranylcypromine-propionyl		189	C12H15NO	1000123-86-3	12
5	Cyclopropanamine, 1-phenyl-		133	C9H11N	1000351-26-2	10



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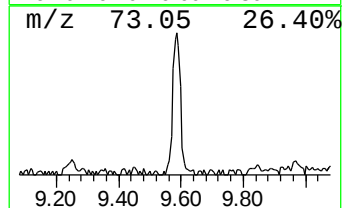
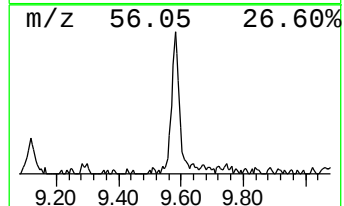
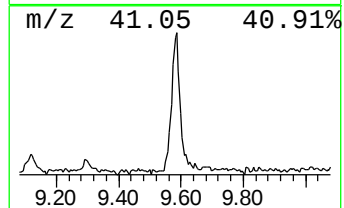
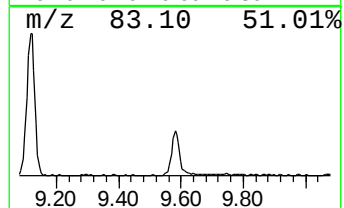
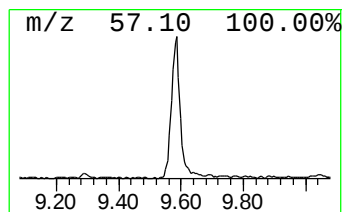
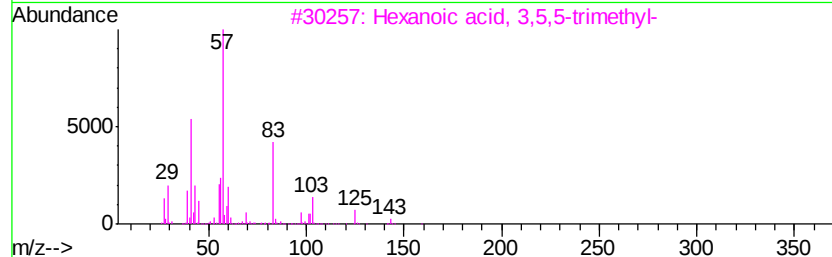
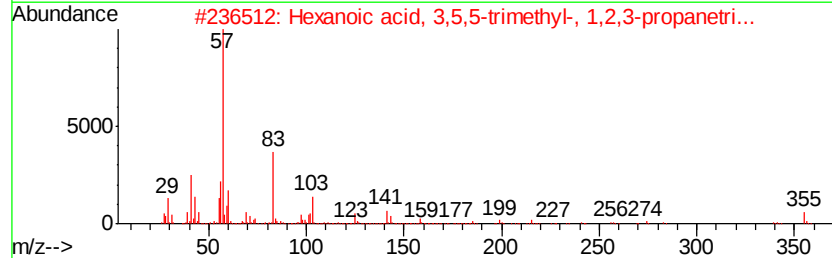
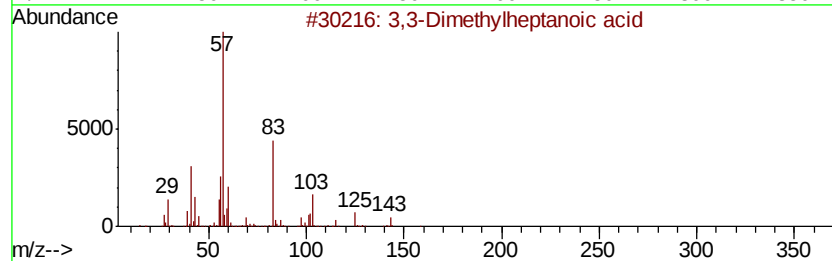
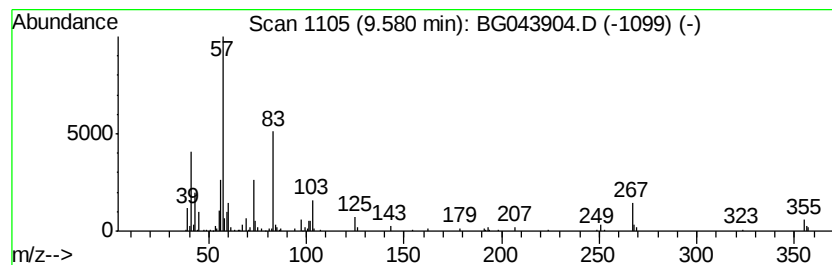
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Peak Number 5 3,3-Dimethylheptanoic acid Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.58	3.27 ng	196909	Naphthalene-d8	10.77

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3,3-Dimethylheptanoic acid	158	C9H18O2	067061-30-7	59
2		Hexanoic acid, 3,5,5-trimethyl-,...	512	C30H56O6	056554-53-1	49
3		Hexanoic acid, 3,5,5-trimethyl-	158	C9H18O2	003302-10-1	47
4		Hexanoyl chloride, 3,5,5-trimethyl-	176	C9H17ClO	036727-29-4	43
5		3-Cyclopentylpropionic acid, 2-b...	198	C12H22O2	1000293-36-5	38



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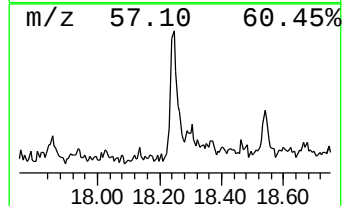
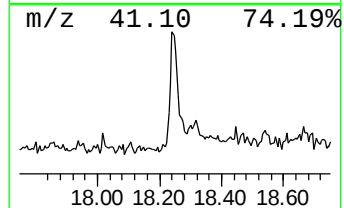
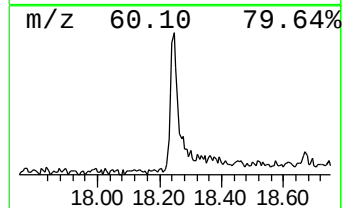
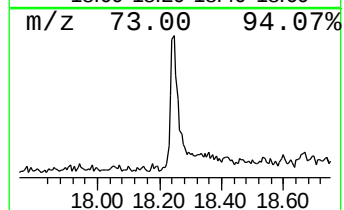
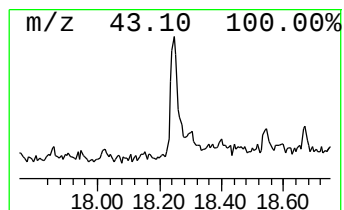
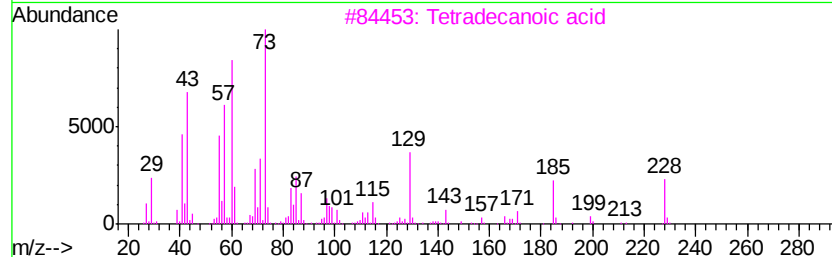
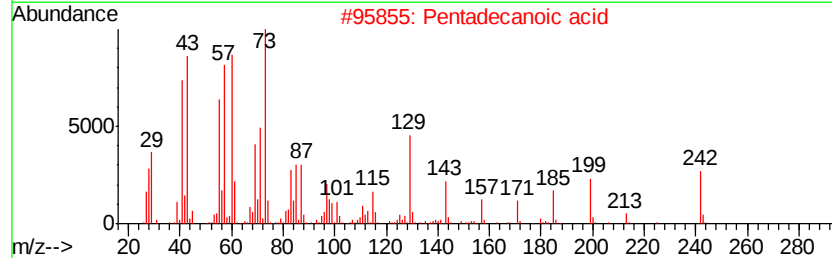
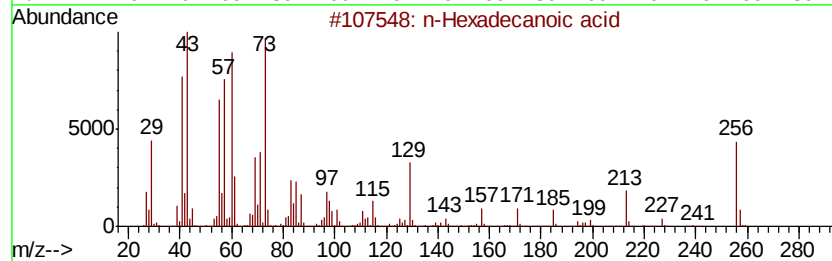
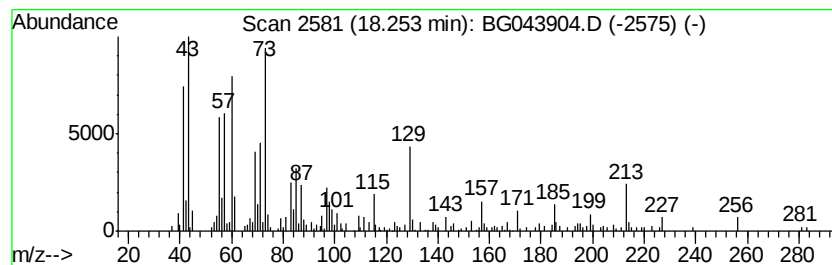
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Peak Number 6 n-Hexadecanoic acid Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD		R.T.
18.25	2.14 ng	228995	Phenanthrene-d10		17.34
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2	Pentadecanoic acid	242	C15H30O2	001002-84-2	87
3	Tetradecanoic acid	228	C14H28O2	000544-63-8	81
4	Octadecanoic acid	284	C18H36O2	000057-11-4	68
5	Tridecanoic acid	214	C13H26O2	000638-53-9	64



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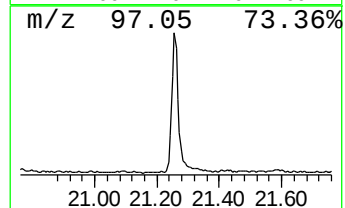
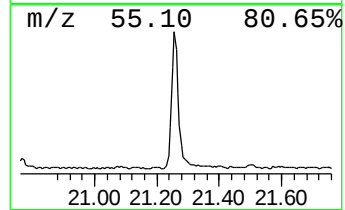
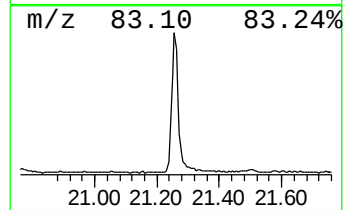
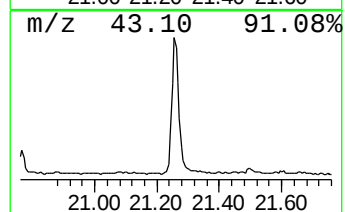
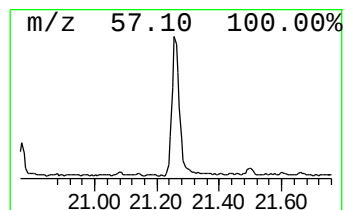
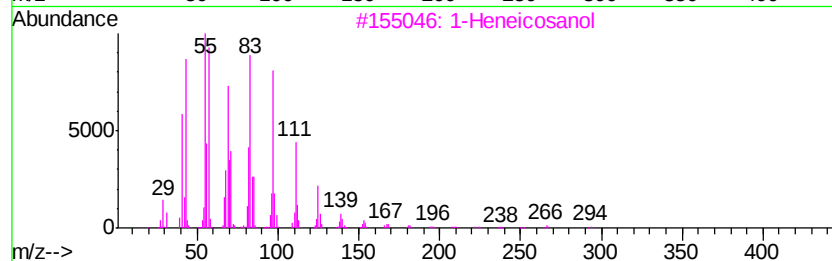
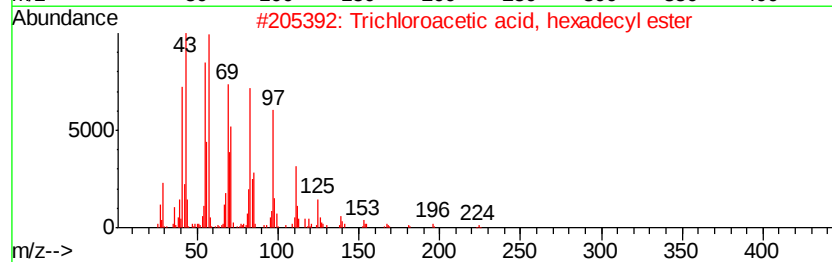
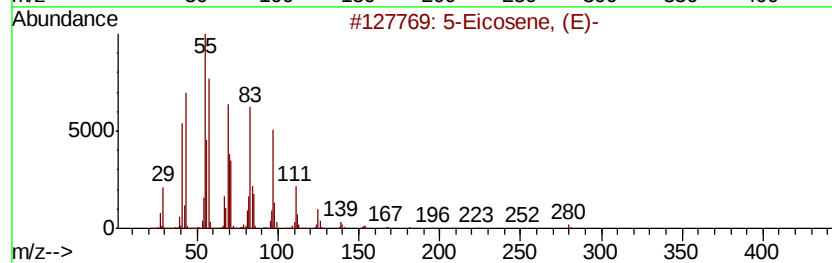
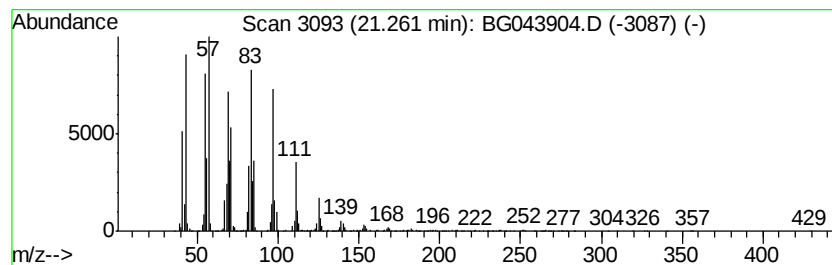
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 7 5-Eicosene, (E)- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.26	16.84 ng	1800500	Chrysene-d12	21.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Eicosene, (E)-	280	C20H40	074685-30-6	99
2		Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	94
3		1-Heneicosanol	312	C21H44O	015594-90-8	94
4		Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	93
5		n-Tetracosanol-1	354	C24H50O	000506-51-4	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010320\
Data File : BG043904.D
Acq On : 4 Jan 2020 1:48
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Sample : K6449-05
Misc :
ALS Vial : 18 Sample Multiplier: 1

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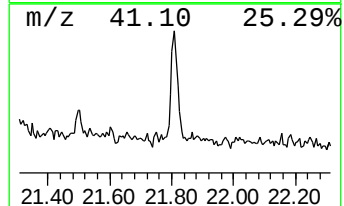
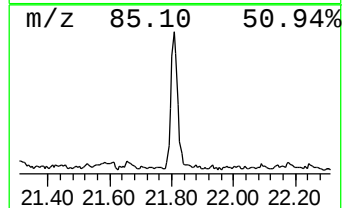
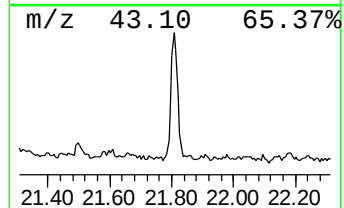
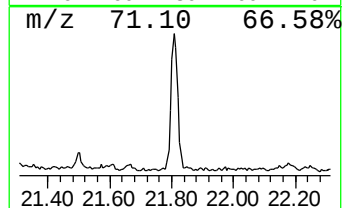
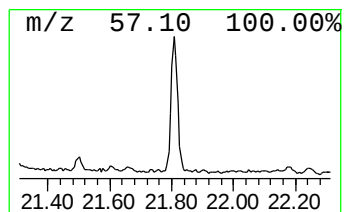
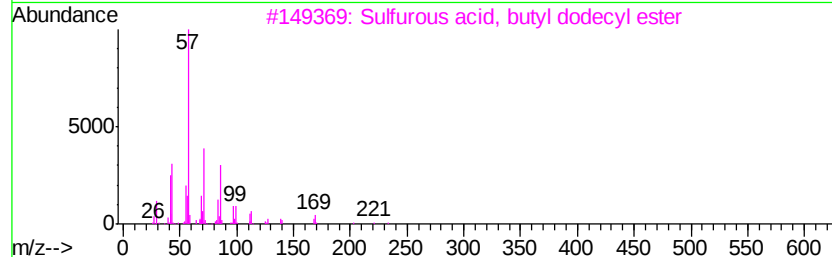
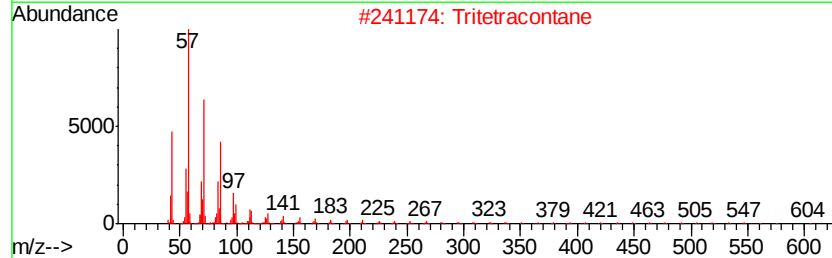
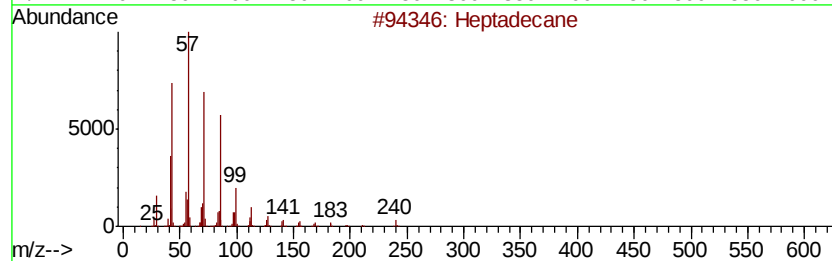
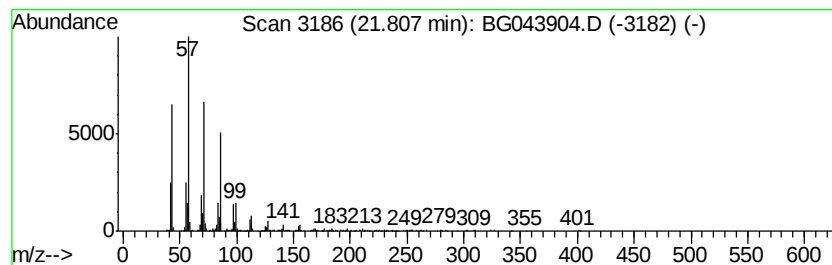
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Heptadecane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.81	3.46 ng	370332	Chrysene-d12	21.58

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane		240	C17H36	000629-78-7	95
2	Tritetracontane		605	C43H88	007098-21-7	91
3	Sulfurous acid, butyl dodecyl ester		306	C16H34O3S	1000309-17-9	91
4	Docosane		310	C22H46	000629-97-0	90
5	Octadecane		254	C18H38	000593-45-3	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010320\
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Misc :
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ClientSampleId :
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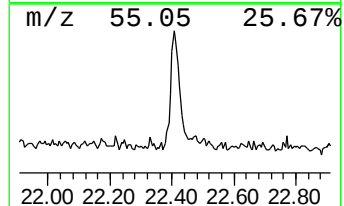
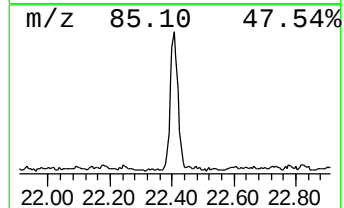
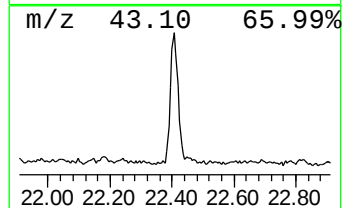
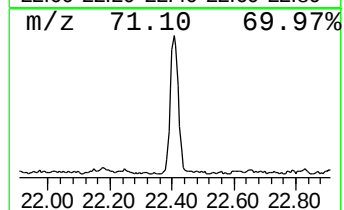
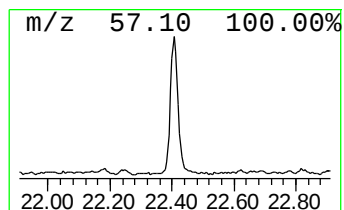
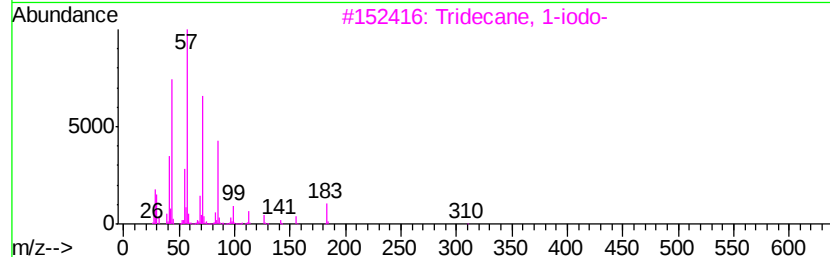
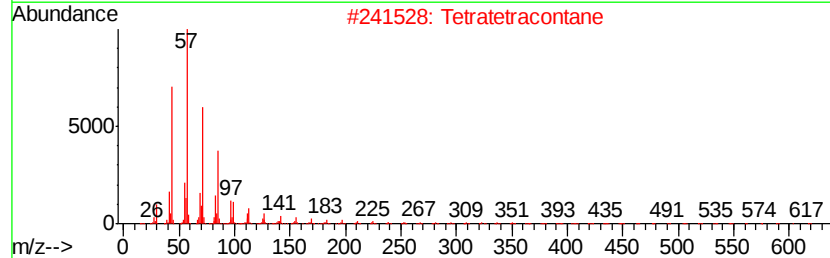
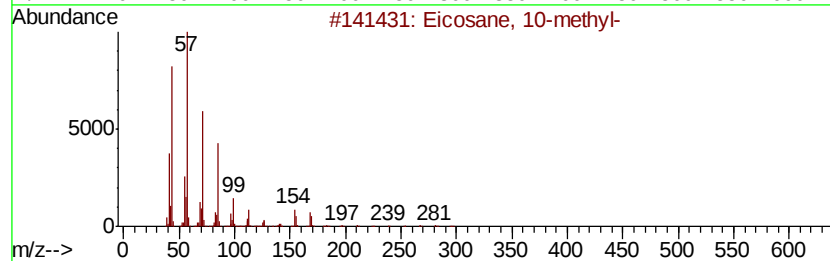
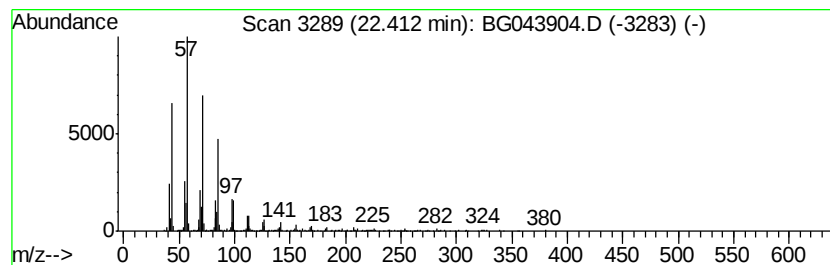
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Eicosane, 10-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.41	5.38 ng	575660	Chrysene-d12	21.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane, 10-methyl-	296	C21H44	054833-23-7	93
2		Tetratetracontane	619	C44H90	007098-22-8	91
3		Tridecane, 1-iodo-	310	C13H27I	035599-77-0	91
4		Octacosane	394	C28H58	000630-02-4	91
5		Tricosane	324	C23H48	000638-67-5	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010320\
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Acq On : 4 Jan 2020 1:48
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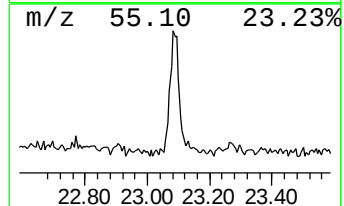
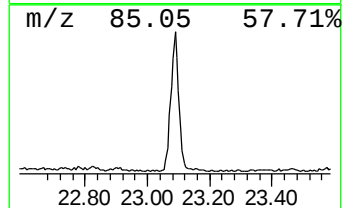
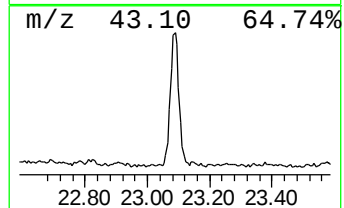
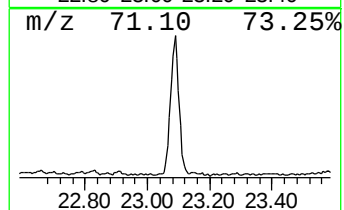
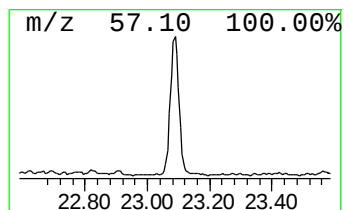
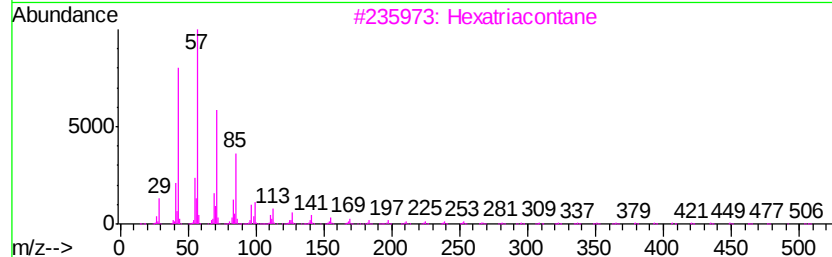
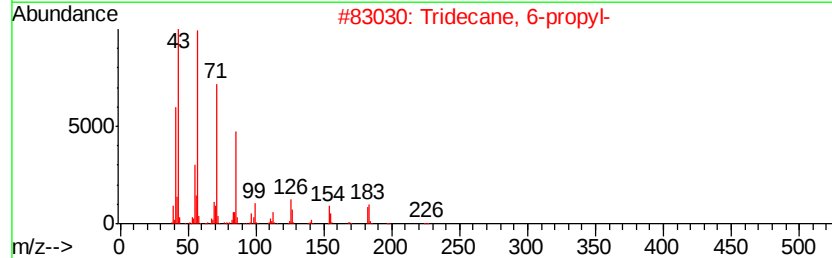
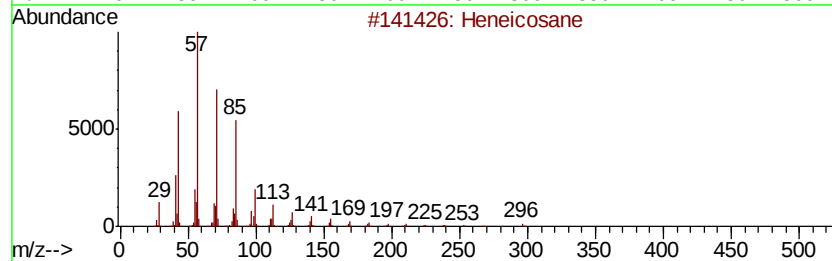
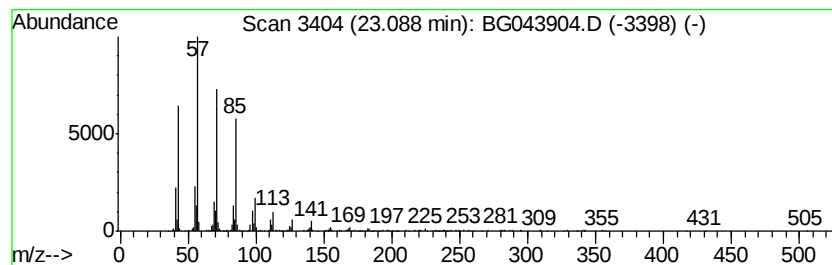
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Heneicosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.09	5.97 ng	638851	Chrysene-d12	21.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	91
2		Tridecane, 6-propyl-	226	C16H34	055045-10-8	87
3		Hexatriacontane	507	C36H74	000630-06-8	80
4		Heptacosane	380	C27H56	000593-49-7	80
5		Tetracosane	338	C24H50	000646-31-1	80



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010320\
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Acq On : 4 Jan 2020 1:48
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ClientSampleId :
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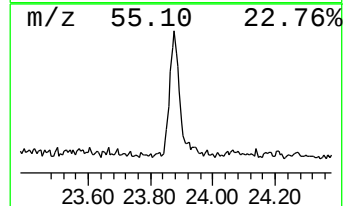
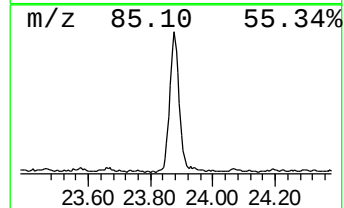
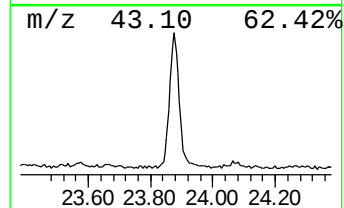
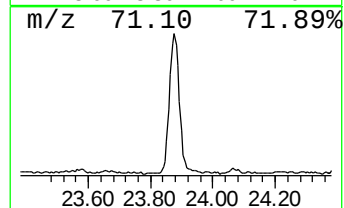
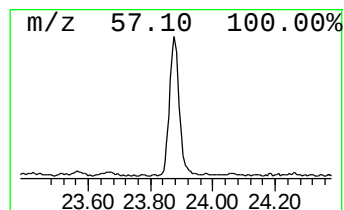
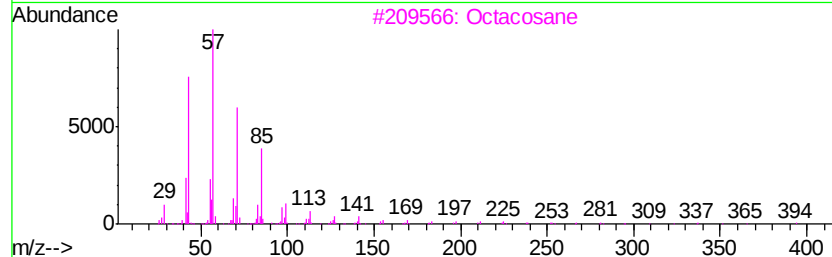
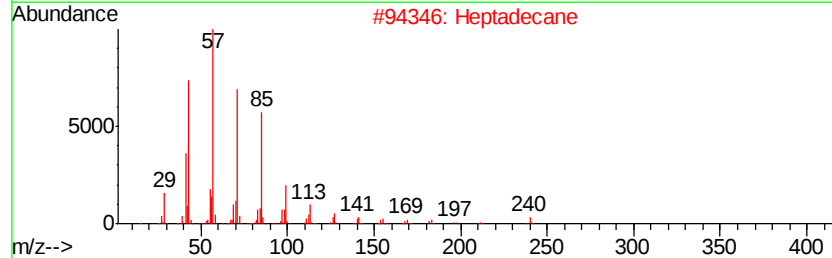
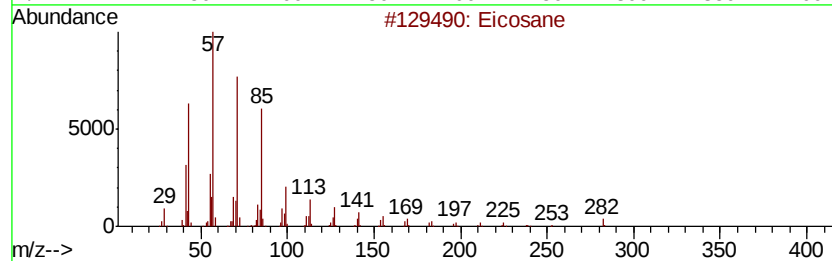
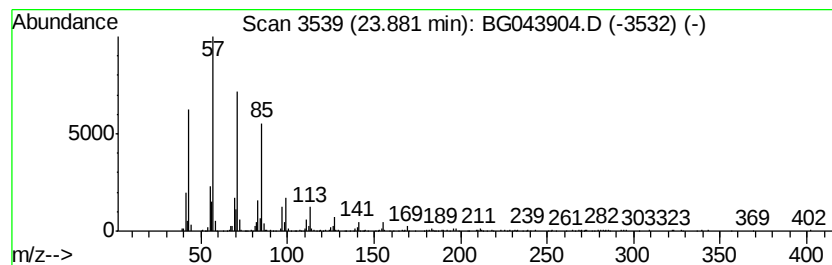
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 Eicosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.88	6.37 ng	693313	Perylene-d12	24.72

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	94
2		Heptadecane	240	C17H36	000629-78-7	93
3		Octacosane	394	C28H58	000630-02-4	90
4		Tetracosane	338	C24H50	000646-31-1	87
5		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010320\
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Acq On : 4 Jan 2020 1:48
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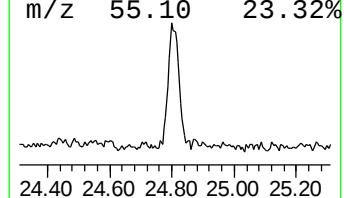
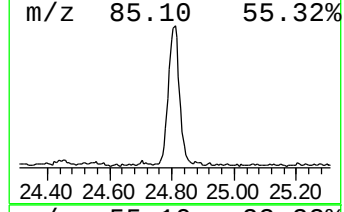
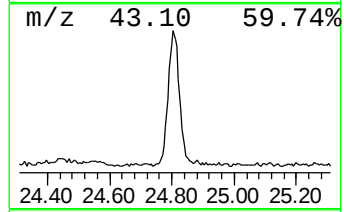
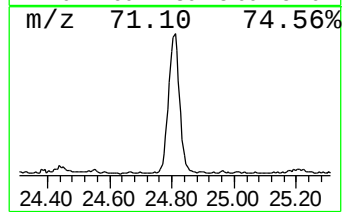
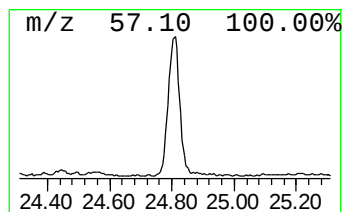
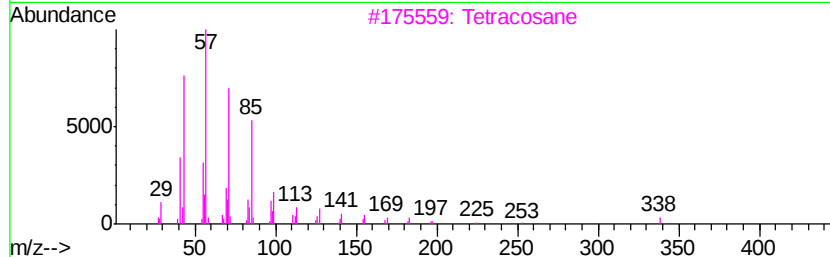
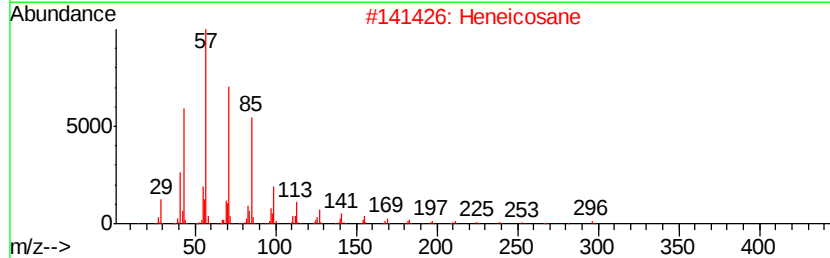
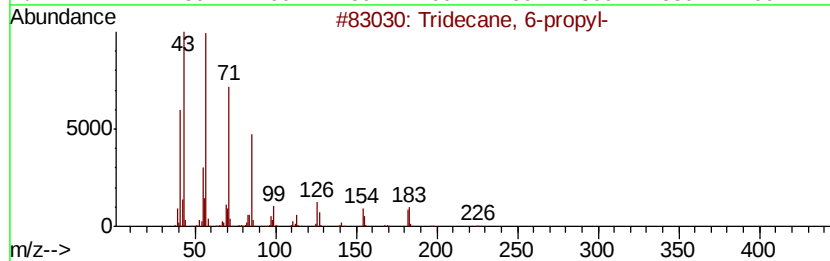
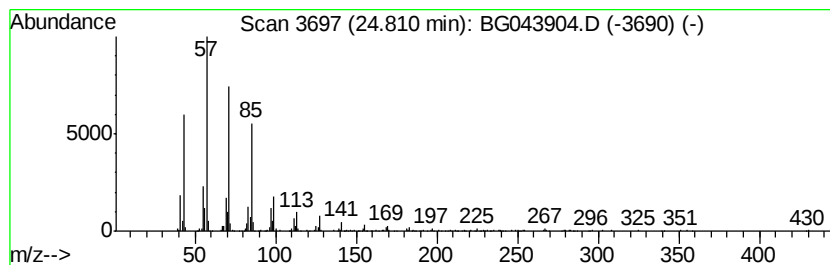
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Tridecane, 6-propyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.81	6.38 ng	695048	Perylene-d12	24.72

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tridecane, 6-propyl-	226	C16H34	055045-10-8	91
2		Heneicosane	296	C21H44	000629-94-7	91
3		Tetracosane	338	C24H50	000646-31-1	91
4		Eicosane, 7-hexyl-	366	C26H54	055333-99-8	90
5		Eicosane	282	C20H42	000112-95-8	90



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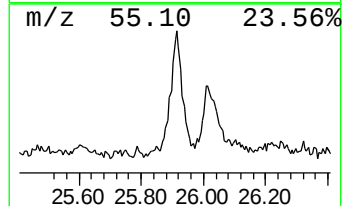
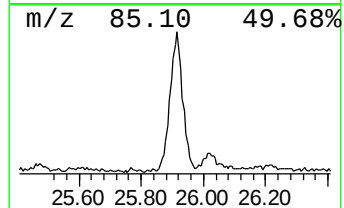
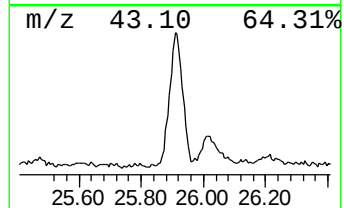
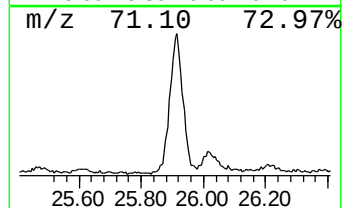
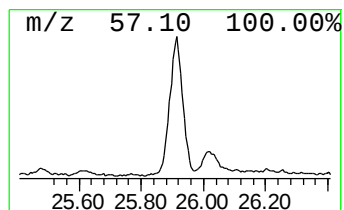
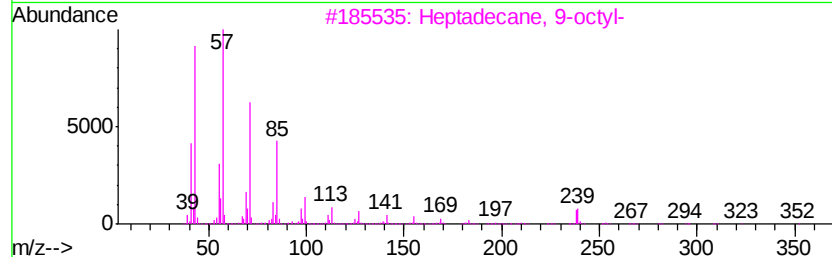
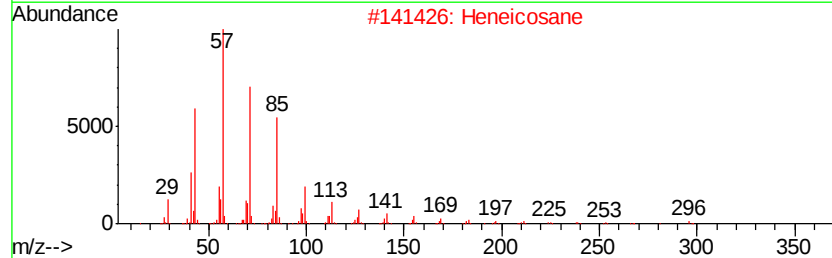
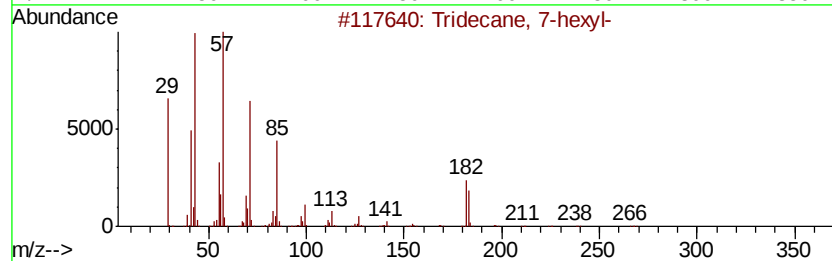
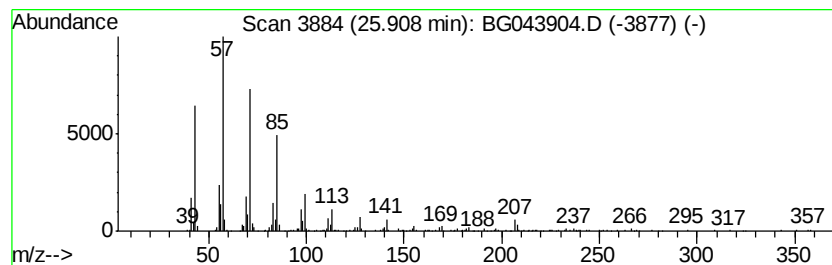
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Tridecane, 7-hexyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.91	5.17 ng	563545	Perylene-d12	24.72

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tridecane, 7-hexyl-	268	C19H40	007225-66-3	93
2		Heneicosane	296	C21H44	000629-94-7	91
3		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
4		Hexacosane	366	C26H54	000630-01-3	90
5		Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG010320\
 Data File : BG043904.D
 Acq On : 4 Jan 2020 1:48
 Operator : JU
 Sample : K6449-05
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
 OW-5

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG123019.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...	3.15	85.2	ng	3369440	1	7.96	791355	20.0
2-Pentanone, 4-hy...	5.07	13.1	ng	517707	1	7.96	791355	20.0
unknown-01	7.50	111.5	ng	4411090	1	7.96	791355	20.0
3,3-Dimethylhepta...	9.58	3.3	ng	196909	2	10.77	1204480	20.0
n-Hexadecanoic acid	18.25	2.1	ng	228995	4	17.34	2144240	20.0
5-Eicosene, (E)-	21.26	16.8	ng	1800500	5	21.58	2138690	20.0
Heptadecane	21.81	3.5	ng	370332	5	21.58	2138690	20.0
Eicosane, 10-methyl-	22.41	5.4	ng	575660	5	21.58	2138690	20.0
Heneicosane	23.09	6.0	ng	638851	5	21.58	2138690	20.0
Eicosane	23.88	6.4	ng	693313	6	24.72	2178170	20.0
Tridecane, 6-propyl-	24.81	6.4	ng	695048	6	24.72	2178170	20.0
Tridecane, 7-hexyl-	25.91	5.2	ng	563545	6	24.72	2178170	20.0