

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021219\
 Data File : BG039464.D
 Acq On : 12 Feb 2019 15:13
 Operator : JU/SJ
 Sample : K1462-11
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 TP-2-S

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-BG012919.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	3.222	17	23	32	rBV	73931	146825	3.47%	0.545%
2	3.292	32	35	44	rVB	52400	78004	1.84%	0.290%
3	5.254	363	369	378	rBB	68533	111558	2.64%	0.414%
4	5.713	436	447	458	rBV	1154632	1885536	44.56%	7.000%
5	7.310	709	719	735	rBV	1279054	2248741	53.14%	8.348%
6	7.698	776	785	793	rBV	1138707	1990520	47.04%	7.390%
7	8.174	857	866	876	rVB	184446	321984	7.61%	1.195%
8	8.497	912	921	929	rBV	766962	1375623	32.51%	5.107%
9	9.349	1056	1066	1074	rBV	708405	1301586	30.76%	4.832%
10	10.994	1338	1346	1354	rVB	286123	511237	12.08%	1.898%
11	13.420	1751	1759	1773	rBV	2007324	3156770	74.60%	11.719%
12	14.236	1892	1898	1904	rBV	145900	210826	4.98%	0.783%
13	14.794	1986	1993	2003	rBV2	511264	816981	19.31%	3.033%
14	16.281	2239	2246	2261	rBV	1646903	2598019	61.39%	9.645%
15	17.538	2453	2460	2467	rBV	661707	1012730	23.93%	3.760%
16	18.395	2597	2606	2623	rBV2	153470	281122	6.64%	1.044%
17	18.730	2657	2663	2668	rBV7	44245	63835	1.51%	0.237%
18	19.653	2816	2820	2833	rBV	81516	146769	3.47%	0.545%
19	19.935	2861	2868	2875	rVB8	22942	45778	1.08%	0.170%
20	20.134	2895	2902	2907	rBV	3123938	4231720	100.00%	15.710%
21	20.763	3004	3009	3013	rBV3	37051	54089	1.28%	0.201%
22	20.904	3029	3033	3039	rVB	41914	49363	1.17%	0.183%
23	21.421	3116	3121	3133	rBV	169359	271374	6.41%	1.007%
24	21.826	3183	3190	3201	rVB	634494	1083364	25.60%	4.022%
25	21.985	3213	3217	3222	rVB	75141	105027	2.48%	0.390%
26	22.619	3319	3325	3333	rBV	126555	212414	5.02%	0.789%
27	23.342	3440	3448	3456	rBV	145789	285971	6.76%	1.062%
28	23.547	3477	3483	3489	rBV2	70462	134076	3.17%	0.498%
29	24.188	3584	3592	3600	rBV2	147771	328123	7.75%	1.218%
30	24.270	3600	3606	3613	rVV7	28232	72549	1.71%	0.269%
31	25.204	3751	3765	3778	rBV3	413997	1354872	32.02%	5.030%
32	26.379	3956	3965	3975	rVB	91231	249335	5.89%	0.926%
33	27.812	4200	4209	4221	rVB4	36348	129247	3.05%	0.480%
34	29.551	4497	4505	4518	rVB7	17424	70115	1.66%	0.260%

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

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

Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG012919.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

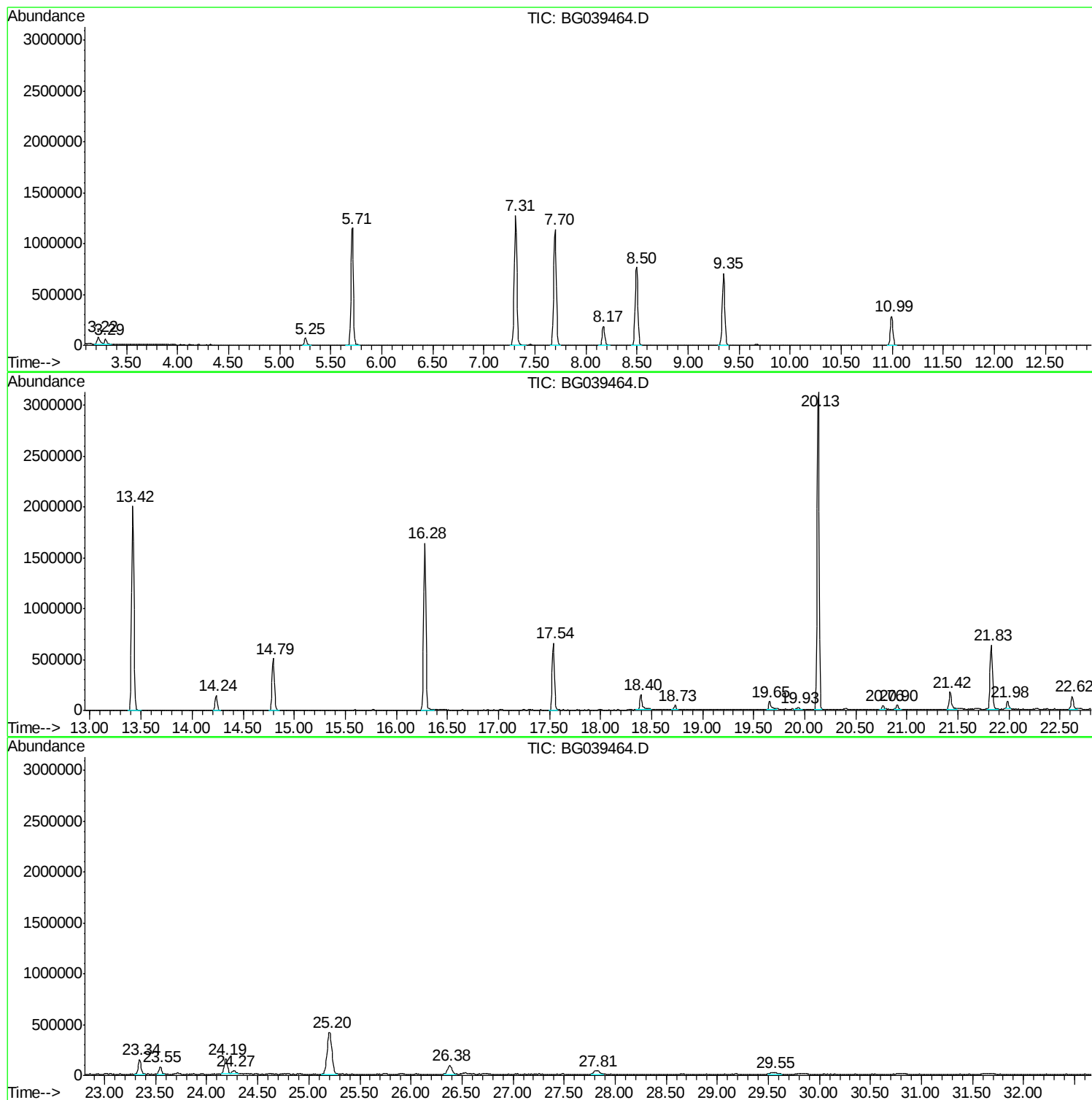
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG012919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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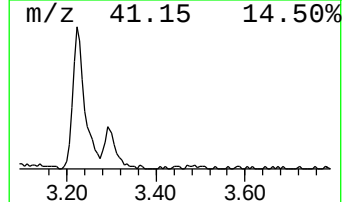
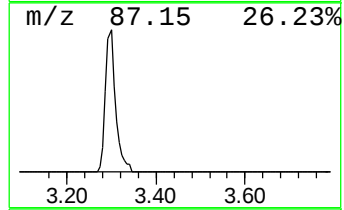
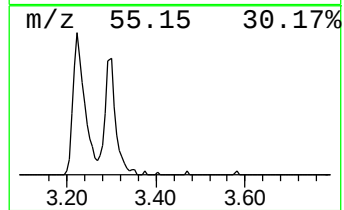
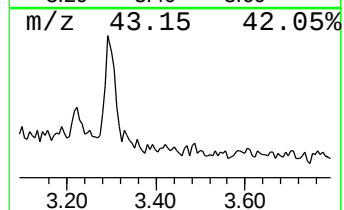
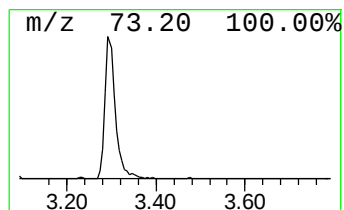
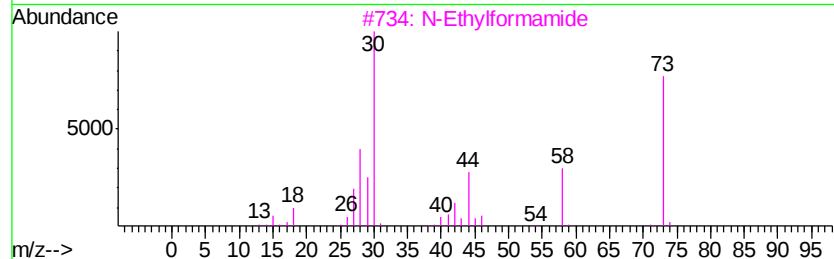
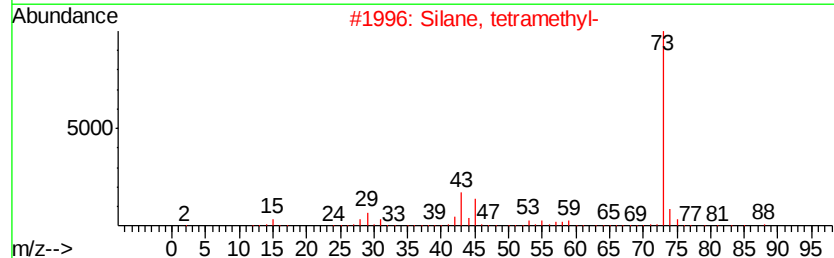
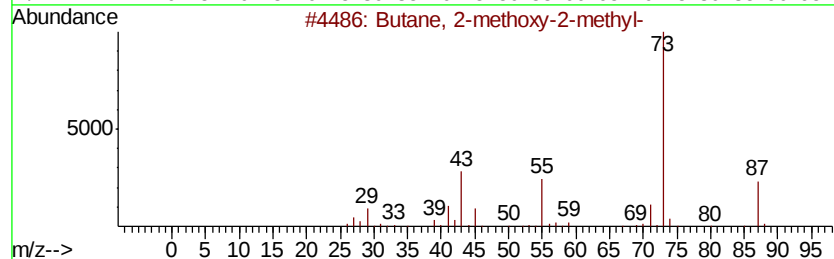
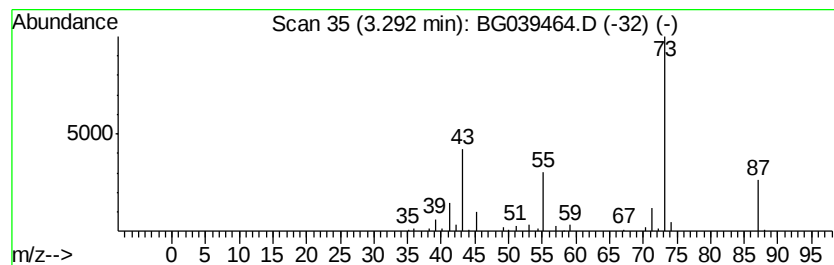
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Butane, 2-methoxy-2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.29	4.85 ng	78004	1,4-Dichlorobenzene-d4	8.17

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2			Silane, tetramethyl-	88	C4H12Si	000075-76-3	10
3			N-Ethylformamide	73	C3H7NO	000627-45-2	9
4			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9
5			3-Pentanol, 3-methyl-	102	C6H14O	000077-74-7	9



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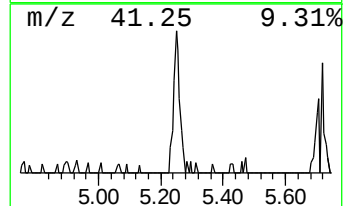
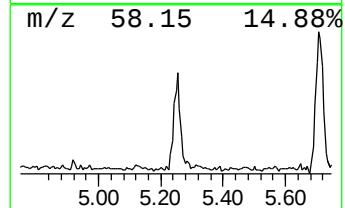
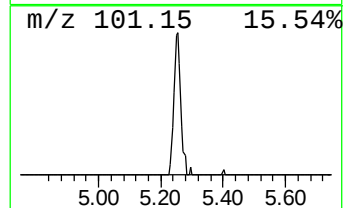
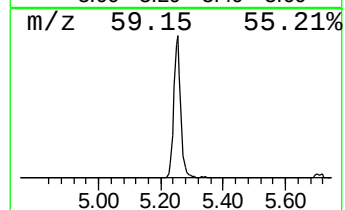
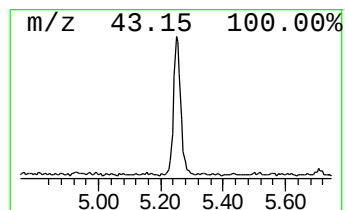
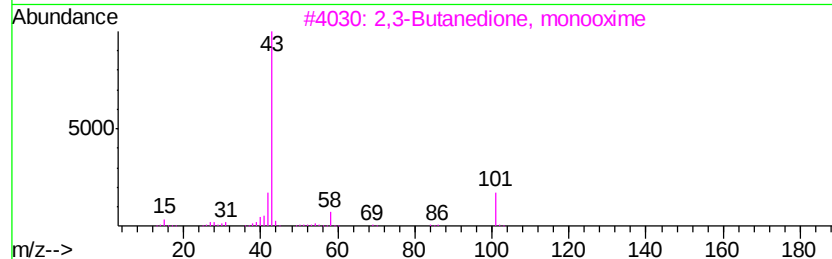
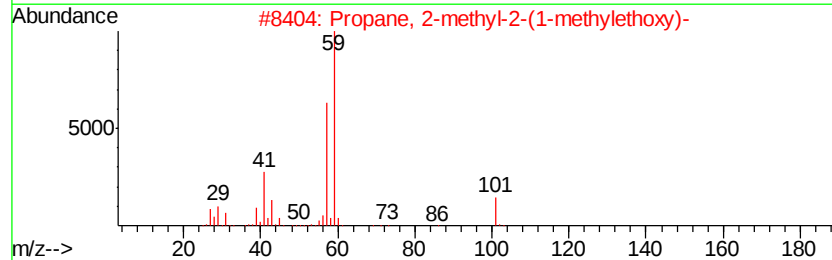
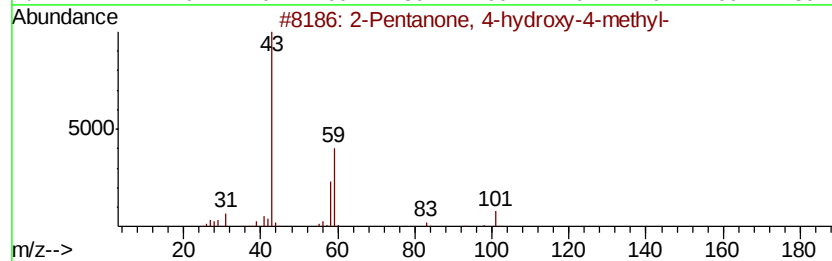
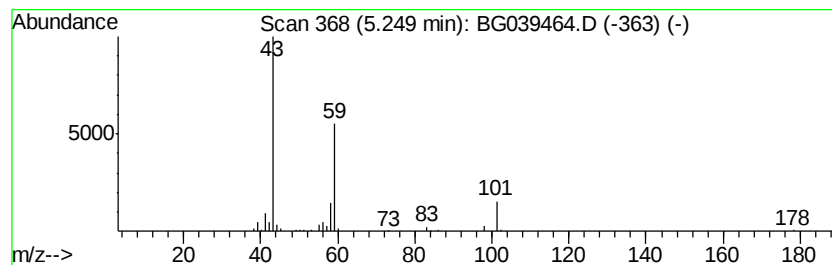
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.25	6.93 ng	111558	1,4-Dichlorobenzene-d4	8.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2		Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	40
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	35
4		Hydrazine, 1,1-bis(1-methylethyl)-	116	C6H16N2	000921-14-2	28
5		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28



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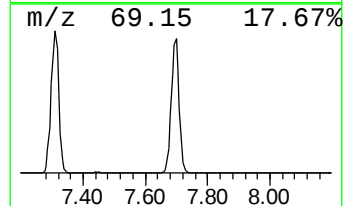
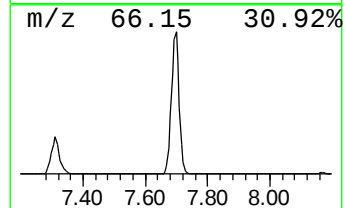
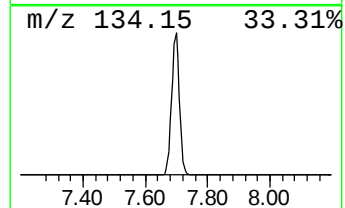
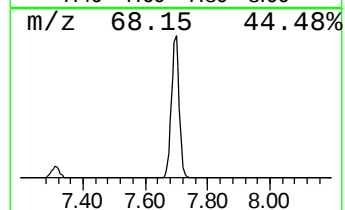
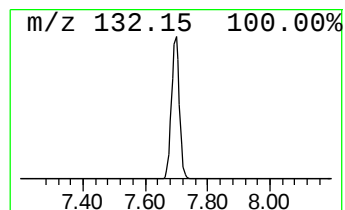
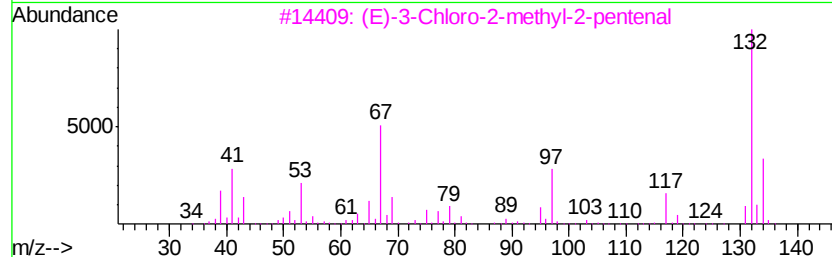
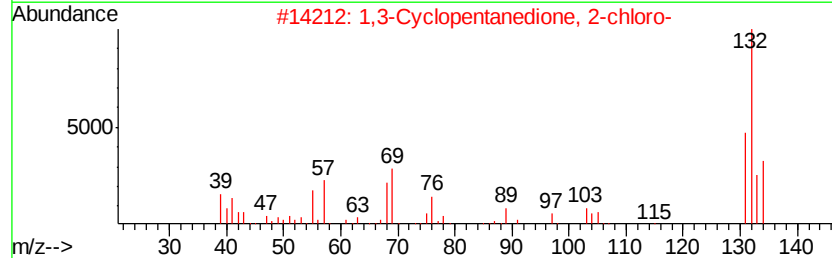
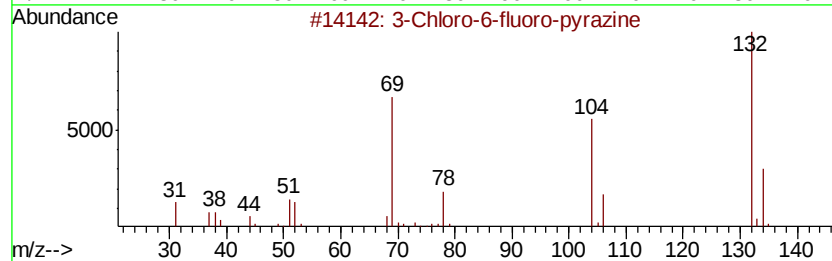
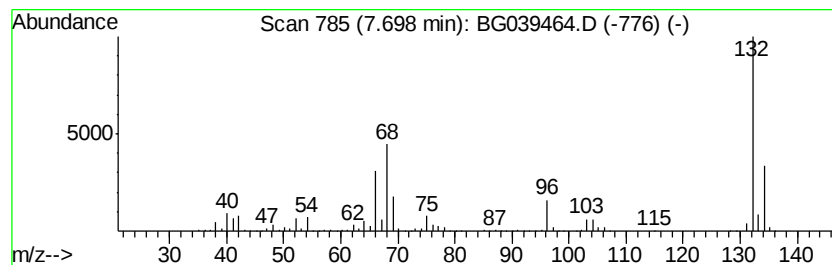
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 unknown7.70 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.70	123.64 ng	1990520	1,4-Dichlorobenzene-d4	8.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	17
3		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
4		(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	10
5		Xenon	132	Xe	007440-63-3	9



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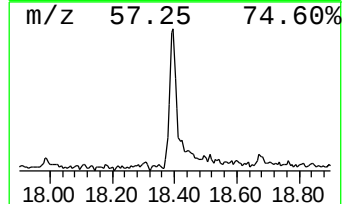
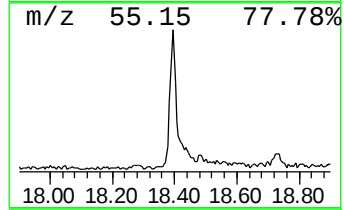
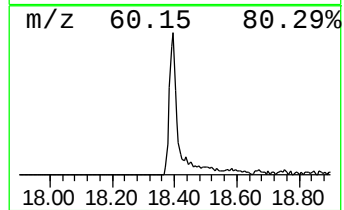
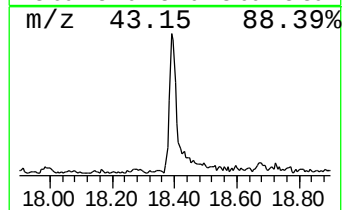
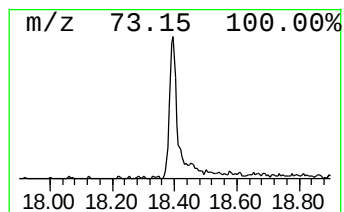
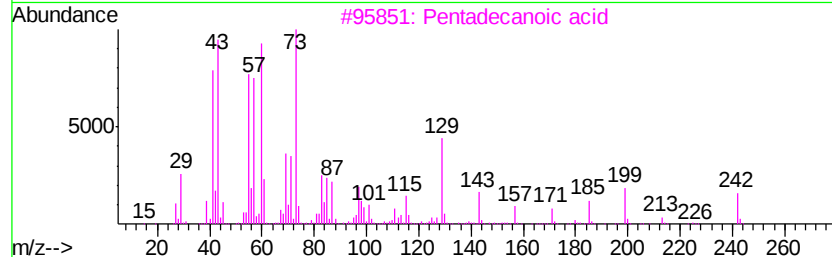
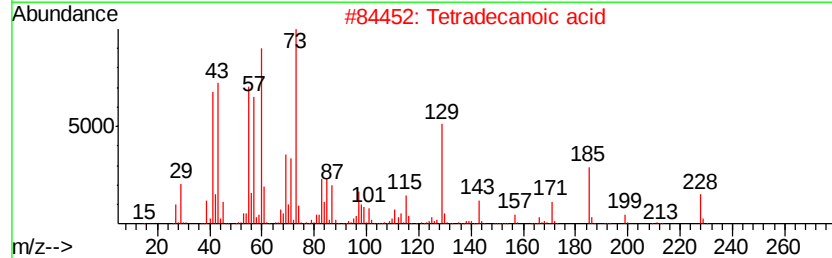
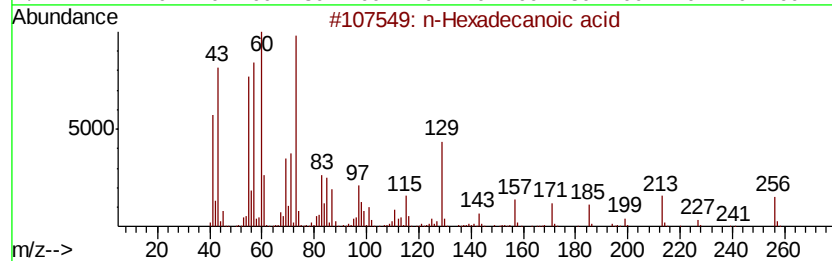
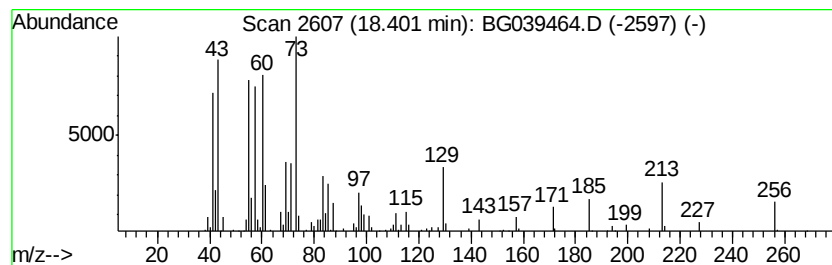
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 n-Hexadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD		R.T.	
18.40	5.55 ng	281122	Phenanthrene-d10		17.54	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		Tetradecanoic acid	228	C14H28O2	000544-63-8	93
3		Pentadecanoic acid	242	C15H30O2	001002-84-2	91
4		Tridecanoic acid	214	C13H26O2	000638-53-9	76
5		Dodecanoic acid	200	C12H24O2	000143-07-7	64



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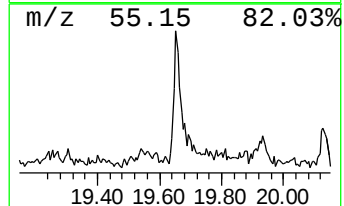
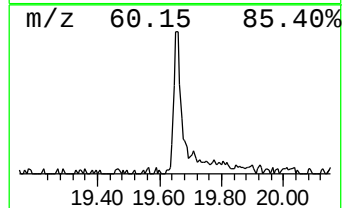
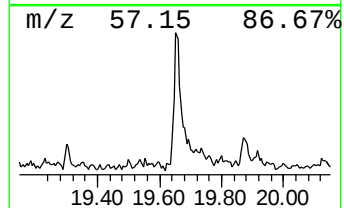
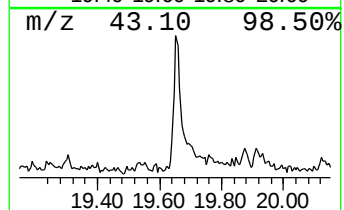
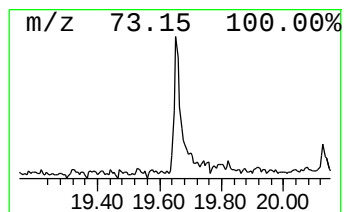
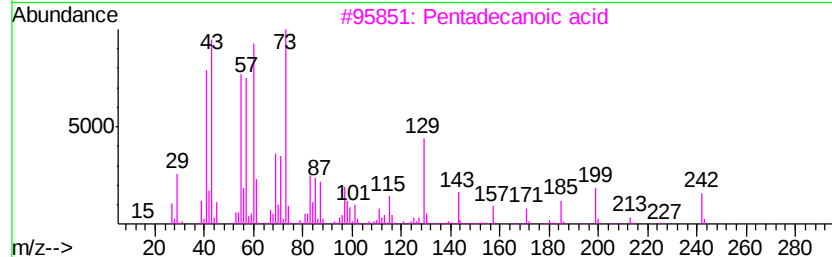
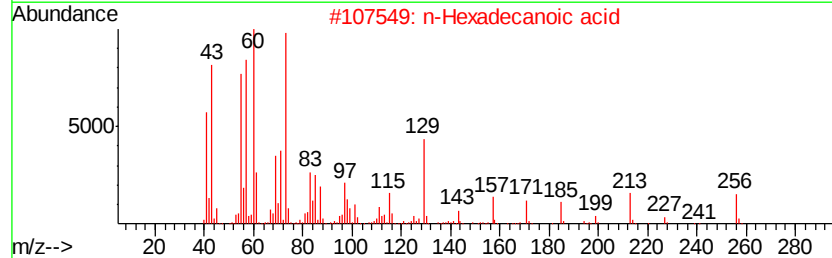
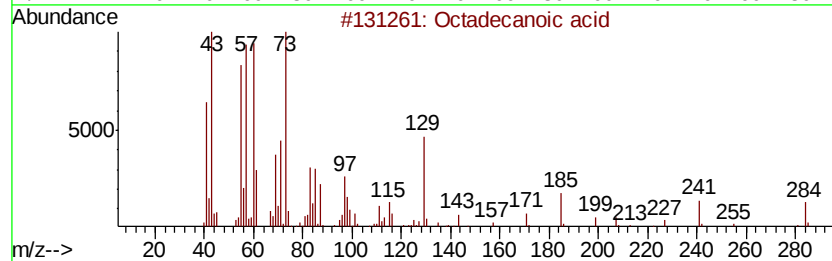
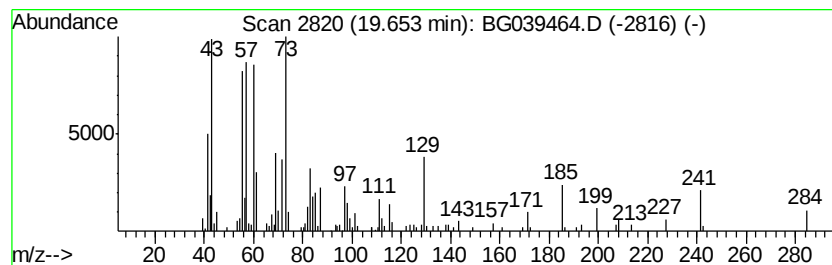
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TIC Library : C:\Database\NIST11.L
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Peak Number 7 Octadecanoic acid Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD		R.T.
19.65	2.90 ng	146769	Phenanthrene-d10		17.54
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	98
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	81
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	72
4	Tetradecanoic acid	228	C14H28O2	000544-63-8	64
5	n-Decanoic acid	172	C10H20O2	000334-48-5	58



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 Operator : JU/SJ
 Sample : K1462-11
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 TP-2-S

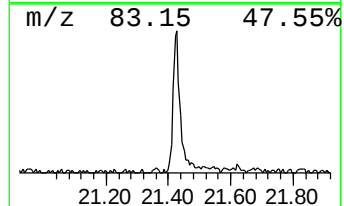
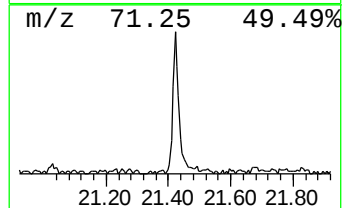
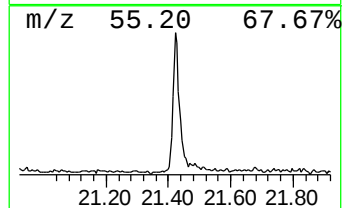
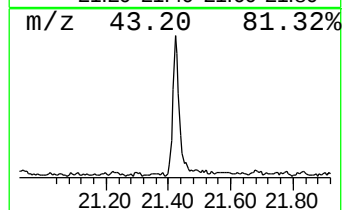
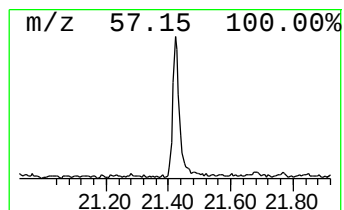
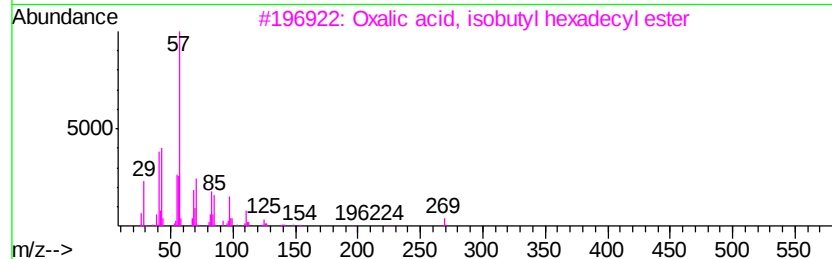
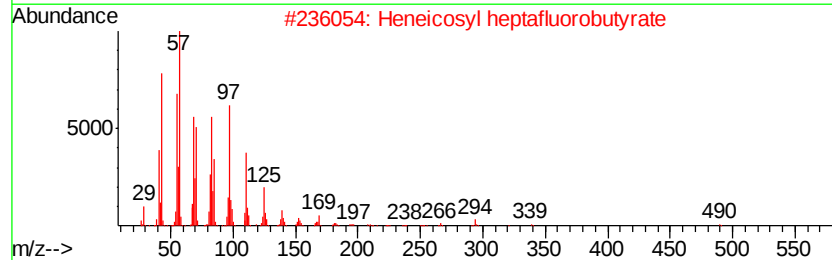
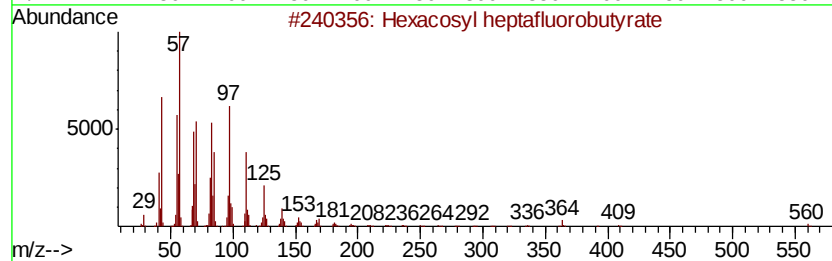
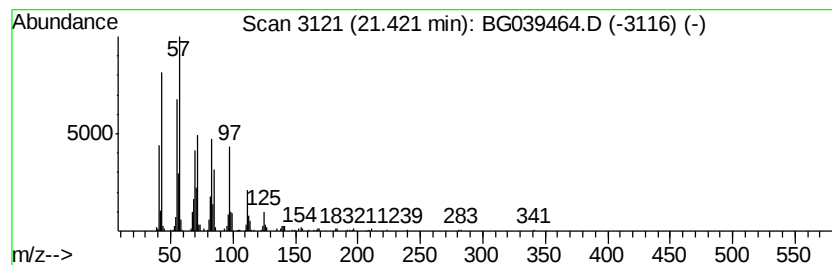
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 8 Hexacosyl heptafluorobutyrate Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.42	5.01 ng	271374	Chrysene-d12	21.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexacosyl heptafluorobutyrate	578	C30H53F7O2	1000351-83-3	91
2		Heneicosyl heptafluorobutyrate	508	C25H43F7O2	1000351-83-8	91
3		Oxalic acid, isobutyl hexadecyl ...	370	C22H42O4	1000309-38-1	91
4		Tricosyl pentafluoropropionate	486	C26H47F5O2	1000351-81-0	91
5		Octacosyl trifluoroacetate	506	C30H57F3O2	1000351-74-9	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021219\
Data File : BG039464.D
Acq On : 12 Feb 2019 15:13
Operator : JU/SJ
Sample : K1462-11
Misc :
ALS Vial : 7 Sample Multiplier: 1

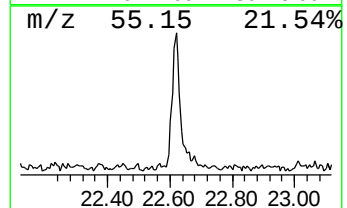
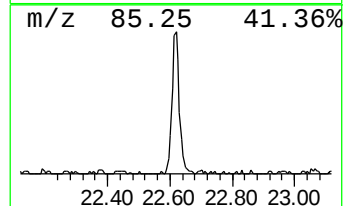
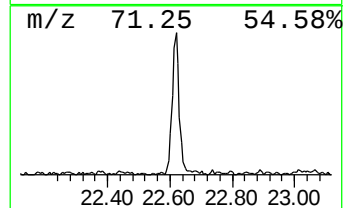
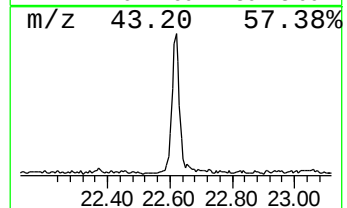
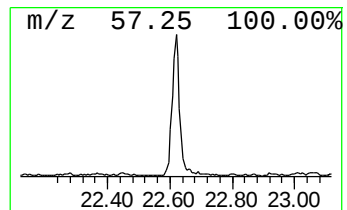
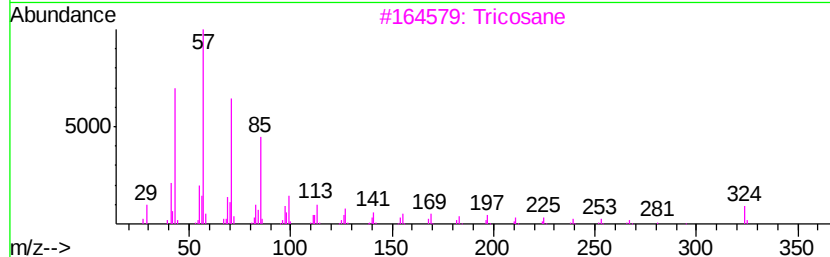
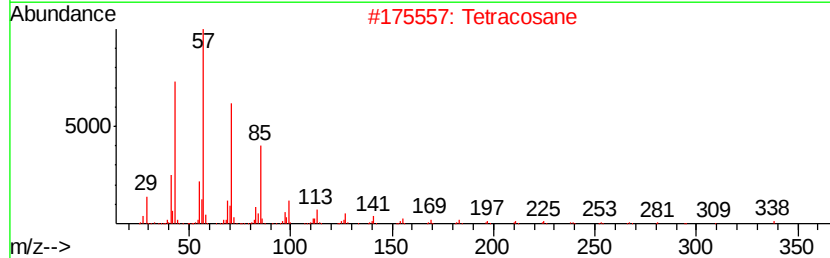
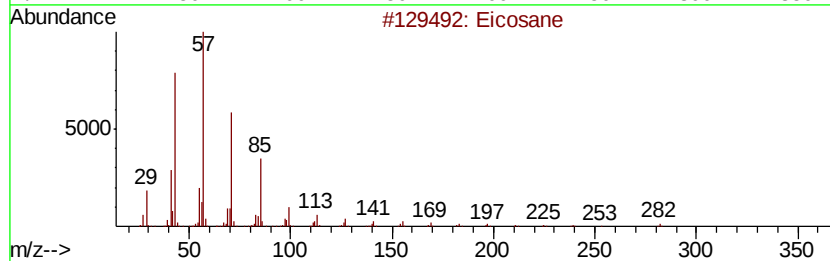
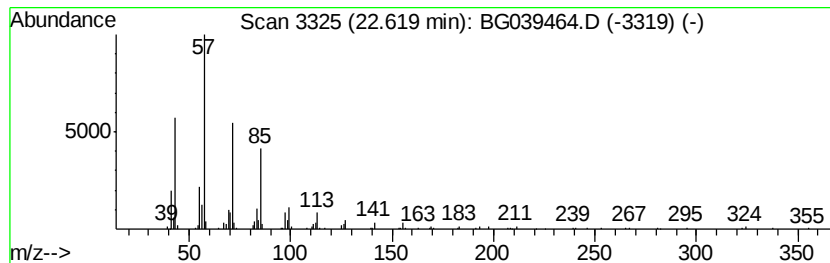
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BNA_G
ClientSampleId :
TP-2-S

Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG012919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Tetracosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.	
22.62	3.92 ng	212414	Chrysene-d12	21.83	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane	282	C20H42	000112-95-8	93
2	Tetracosane	338	C24H50	000646-31-1	91
3	Tricosane	324	C23H48	000638-67-5	91
4	Heneicosane	296	C21H44	000629-94-7	91
5	Heptacosane	380	C27H56	000593-49-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021219\
Data File : BG039464.D
Acq On : 12 Feb 2019 15:13
Operator : JU/SJ
Sample : K1462-11
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
TP-2-S

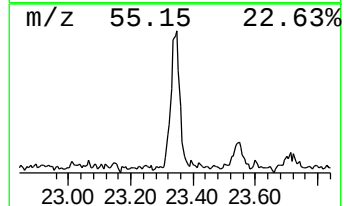
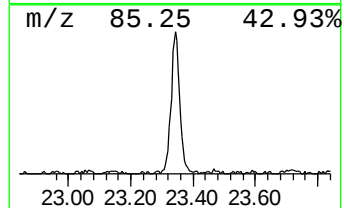
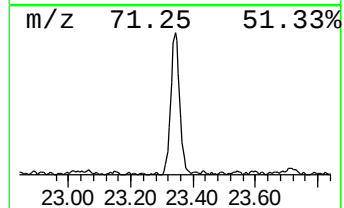
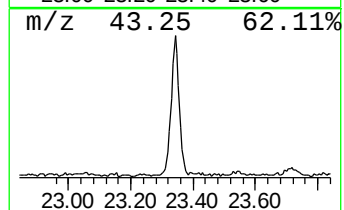
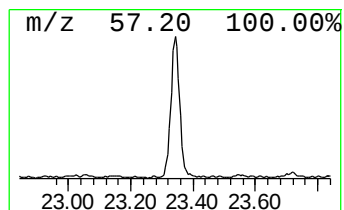
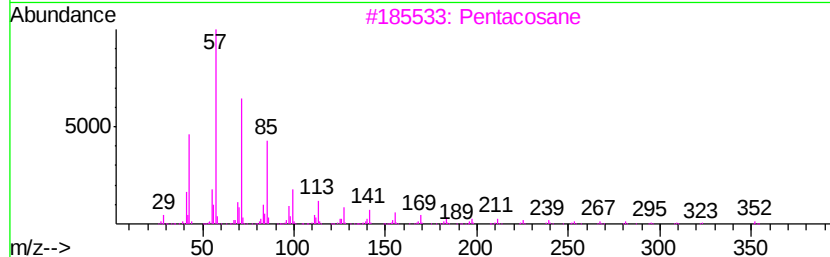
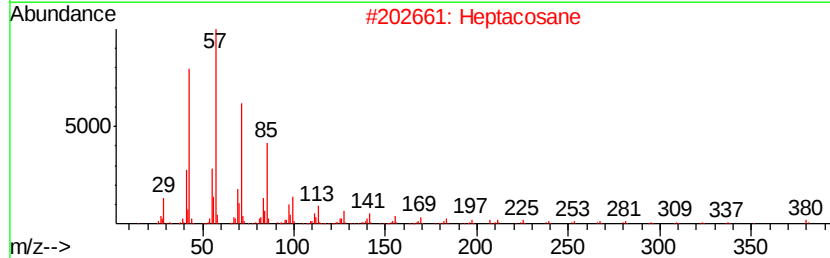
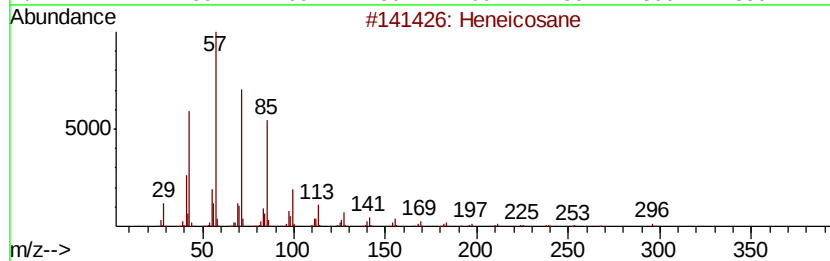
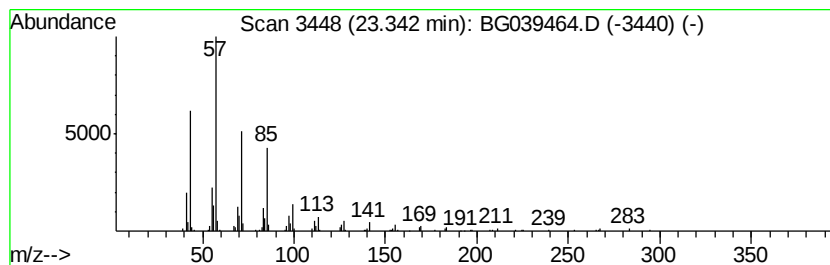
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Heneicosane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.34	5.28 ng	285971	Chrysene-d12	21.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	91
2		Heptacosane	380	C27H56	000593-49-7	91
3		Pentacosane	352	C25H52	000629-99-2	91
4		Tetratetracontane	619	C44H90	007098-22-8	91
5		Hexacosane	366	C26H54	000630-01-3	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021219\
Data File : BG039464.D
Acq On : 12 Feb 2019 15:13
Operator : JU/SJ
Sample : K1462-11
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
TP-2-S

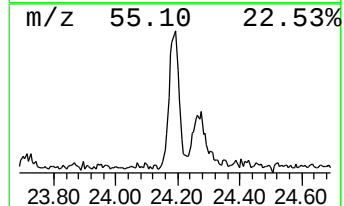
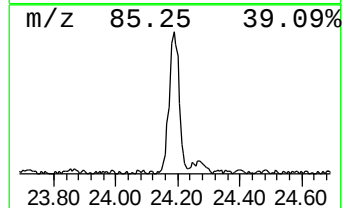
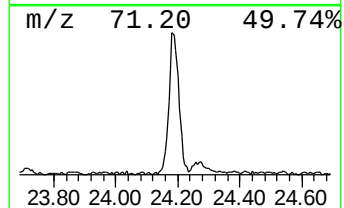
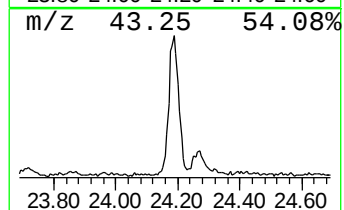
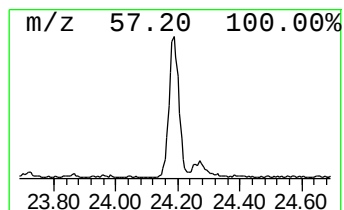
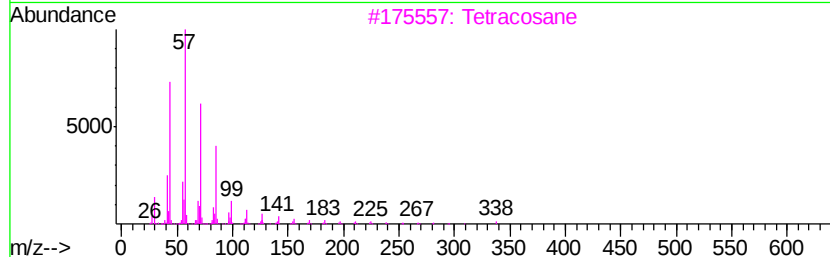
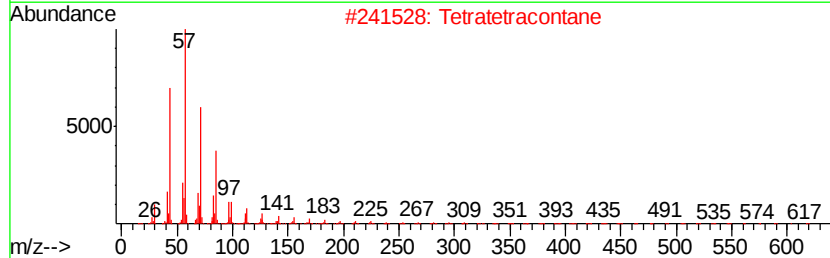
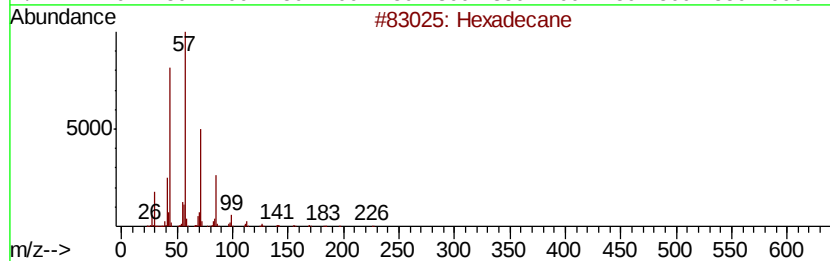
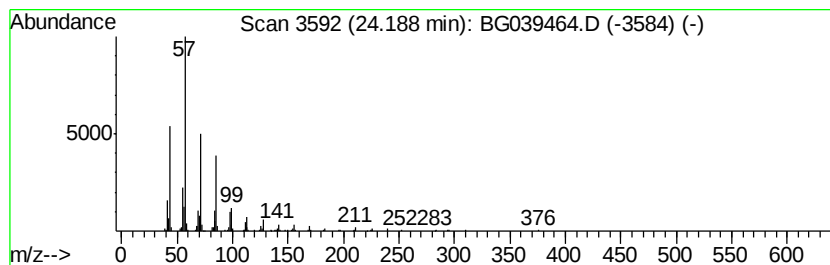
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Hexadecane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.19	4.84 ng	328123	Perylene-d12	25.21

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecane	226	C16H34	000544-76-3	93
2			Tetratetracontane	619	C44H90	007098-22-8	91
3			Tetracosane	338	C24H50	000646-31-1	91
4			2-methyloctacosane	408	C29H60	1000376-72-8	91
5			Octacosane	394	C28H58	000630-02-4	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021219\
Data File : BG039464.D
Acq On : 12 Feb 2019 15:13
Operator : JU/SJ
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Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
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ClientSampleId :
TP-2-S

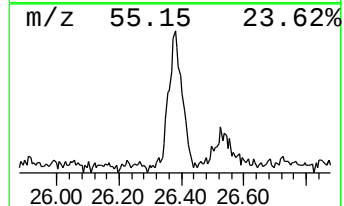
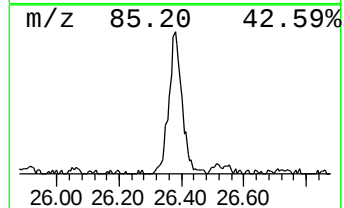
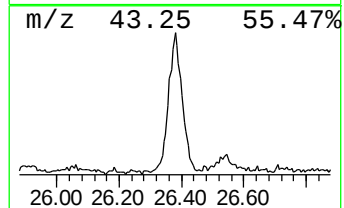
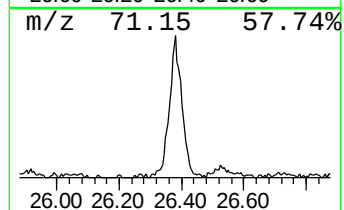
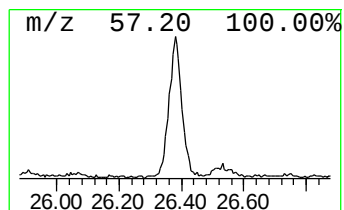
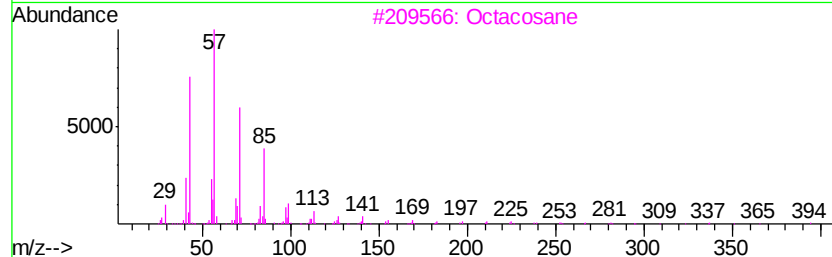
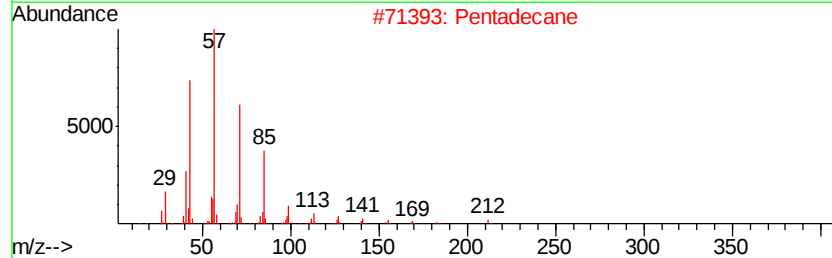
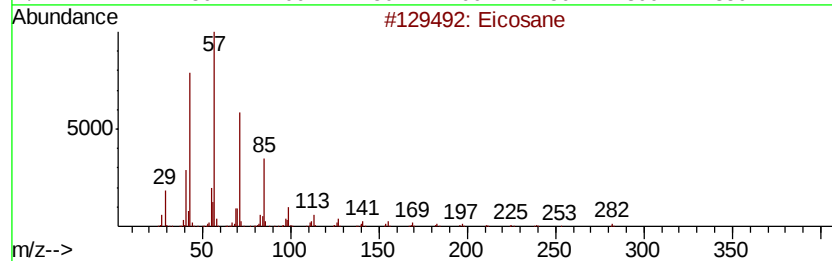
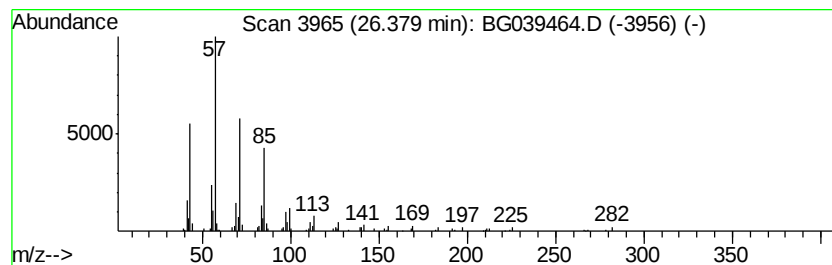
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Eicosane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.38	3.68 ng	249335	Perylene-d12	25.21

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	95
2		Pentadecane	212	C15H32	000629-62-9	91
3		Octacosane	394	C28H58	000630-02-4	90
4		Tetratetracontane	619	C44H90	007098-22-8	90
5		Tetracosane	338	C24H50	000646-31-1	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG021219\
Data File : BG039464.D
Acq On : 12 Feb 2019 15:13
Operator : JU/SJ
Sample : K1462-11
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
TP-2-S

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG012919.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...	3.29	4.8	ng	78004	1	8.17	321984	20.0
2-Pentanone, 4-hy...	5.25	6.9	ng	111558	1	8.17	321984	20.0
unknown7.70	7.70	123.6	ng	1990520	1	8.17	321984	20.0
n-Hexadecanoic acid	18.40	5.5	ng	281122	4	17.54	1012730	20.0
Octadecanoic acid	19.65	2.9	ng	146769	4	17.54	1012730	20.0
Hexacosyl heptafl...	21.42	5.0	ng	271374	5	21.83	1083360	20.0
Tetracosane	22.62	3.9	ng	212414	5	21.83	1083360	20.0
Heneicosane	23.34	5.3	ng	285971	5	21.83	1083360	20.0
Hexadecane	24.19	4.8	ng	328123	6	25.21	1354870	20.0
Eicosane	26.38	3.7	ng	249335	6	25.21	1354870	20.0