

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG103119\
 Data File : BG043439.D
 Acq On : 31 Oct 2019 18:24
 Operator : JU
 Sample : K5587-15
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 CSA-2-1-11

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-BG102119.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	3.196	13	18	27	rVB	125989	229127	9.35%	1.268%
2	5.129	341	347	362	rBV	136983	224029	9.15%	1.240%
3	5.611	422	429	446	rBV	489425	869870	35.51%	4.815%
4	7.209	693	701	712	rBV	506389	986796	40.29%	5.463%
5	7.562	750	761	771	rBV	527243	974859	39.80%	5.397%
6	8.020	832	839	849	rBV	138779	247114	10.09%	1.368%
7	8.349	886	895	904	rBV	414245	757518	30.93%	4.193%
8	9.183	1027	1037	1057	rBV	274923	586665	23.95%	3.248%
9	10.829	1307	1317	1330	rBV	144149	322524	13.17%	1.785%
10	13.273	1726	1733	1747	rBV	835914	1437284	58.68%	7.957%
11	14.101	1865	1874	1890	rBV	100893	194402	7.94%	1.076%
12	14.648	1959	1967	1980	rBV2	310204	516706	21.10%	2.860%
13	16.140	2211	2221	2233	rBV	742679	1301642	53.14%	7.206%
14	17.391	2427	2434	2449	rBV	377067	664709	27.14%	3.680%
15	17.515	2451	2455	2460	rVB4	23313	34134	1.39%	0.189%
16	18.284	2579	2586	2591	rBV5	32552	60795	2.48%	0.337%
17	19.442	2779	2783	2792	rVB2	23681	40640	1.66%	0.225%
18	19.806	2841	2845	2851	rVB	24466	36054	1.47%	0.200%
19	20.000	2872	2878	2894	rBV	1714152	2449304	100.00%	13.559%
20	21.304	3095	3100	3110	rBV4	73090	177191	7.23%	0.981%
21	21.663	3153	3161	3179	rBV2	467595	915615	37.38%	5.069%
22	21.845	3188	3192	3201	rVB3	45968	70513	2.88%	0.390%
23	22.450	3288	3295	3300	rBV3	76280	139959	5.71%	0.775%
24	23.138	3402	3412	3420	rBV4	111701	290726	11.87%	1.609%
25	23.296	3433	3439	3452	rVB4	59080	160960	6.57%	0.891%
26	23.849	3526	3533	3536	rBV4	19016	44813	1.83%	0.248%
27	23.937	3541	3548	3558	rVB3	104905	248640	10.15%	1.376%
28	24.336	3603	3616	3628	rBV3	110867	308985	12.62%	1.710%
29	24.865	3694	3706	3722	rBV3	381640	1114463	45.50%	6.169%
30	25.611	3823	3833	3844	rVB4	146811	480072	19.60%	2.658%
31	25.981	3888	3896	3906	rBV4	58111	167854	6.85%	0.929%
32	27.209	4089	4105	4117	rBV4	171233	703632	28.73%	3.895%
33	27.327	4120	4125	4134	rVB8	29155	82337	3.36%	0.456%
34	29.207	4427	4445	4461	rBV5	163428	849099	34.67%	4.700%

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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	31.733	4859	4875	4876	rBV4	116203	375119	15.32%	2.077%
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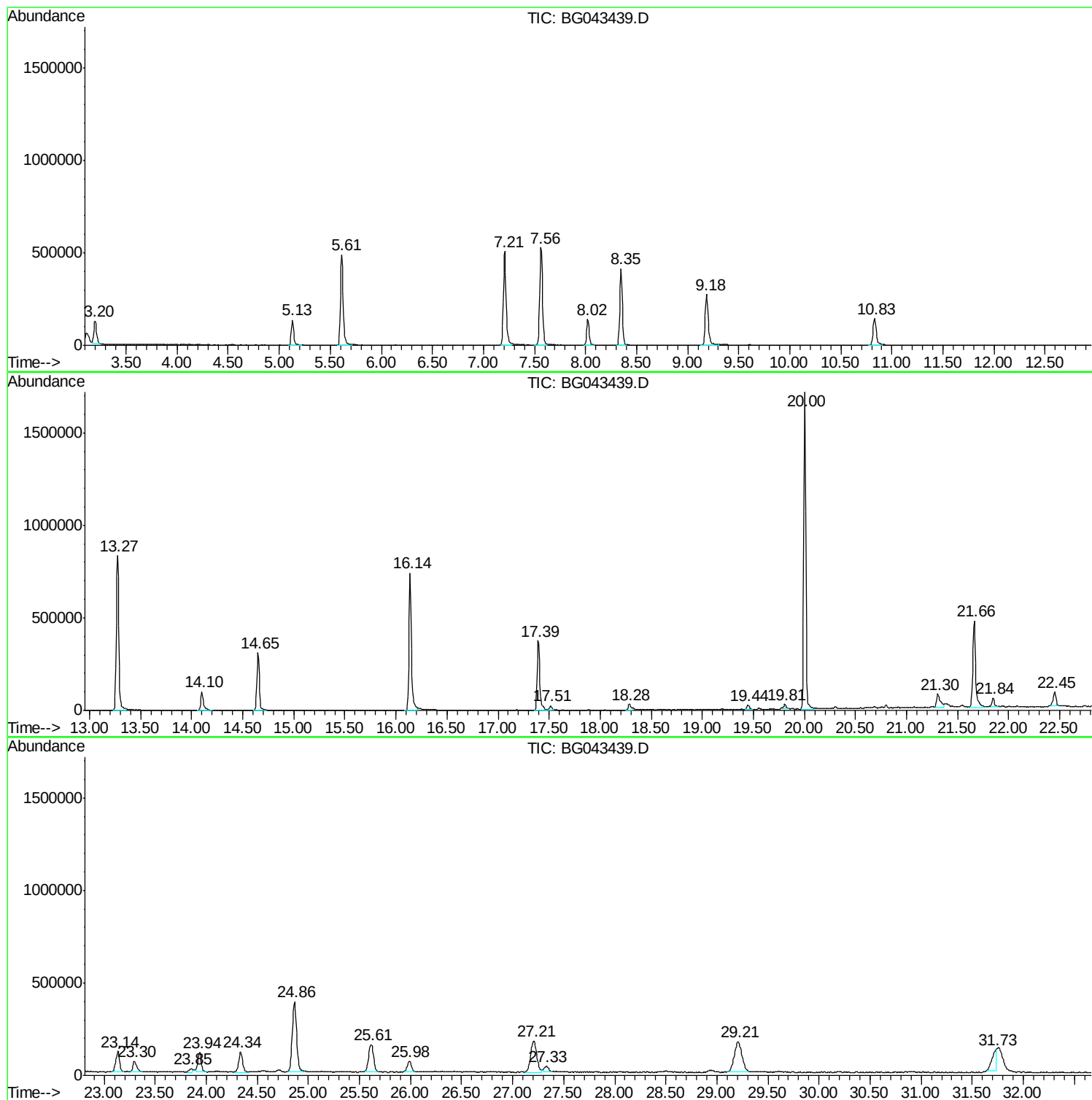
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG102119.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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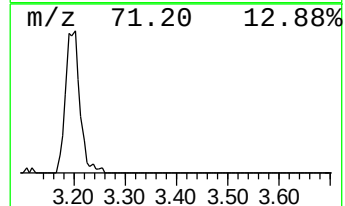
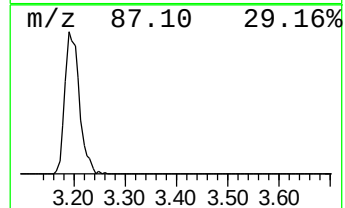
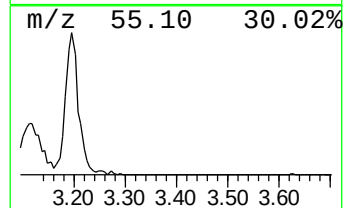
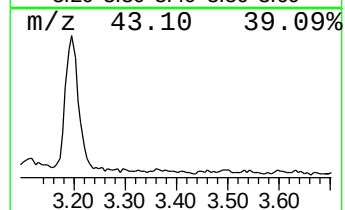
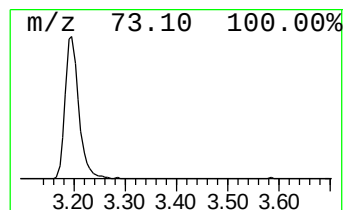
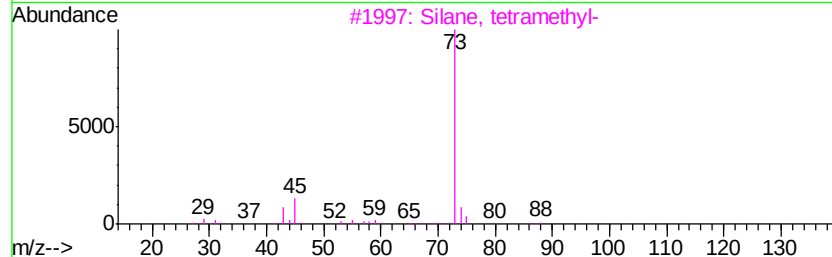
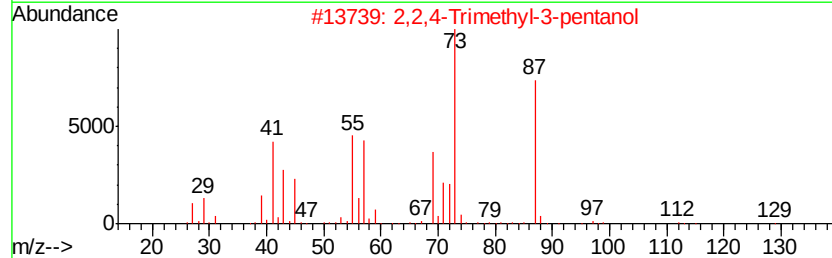
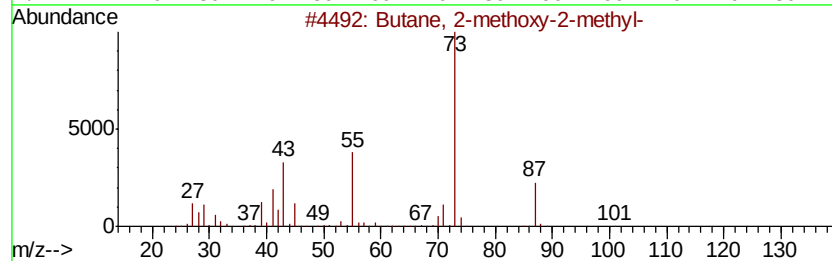
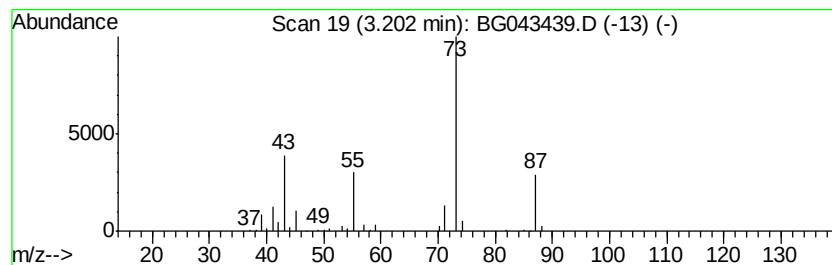
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.20	18.54 ng	229127	1,4-Dichlorobenzene-d4	8.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
3		Silane, tetramethyl-	88	C4H12Si	000075-76-3	9
4		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9
5		N-Ethylformamide	73	C3H7NO	000627-45-2	9



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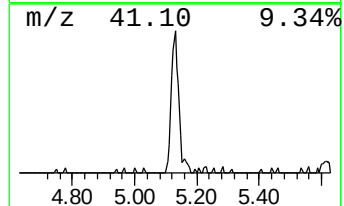
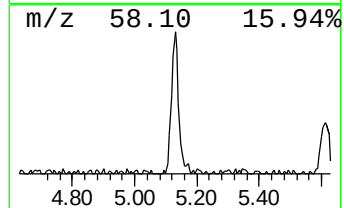
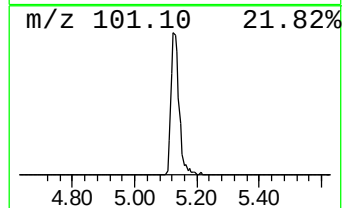
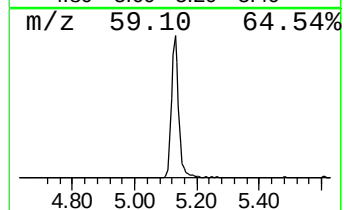
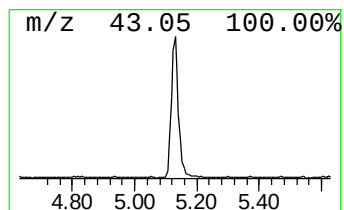
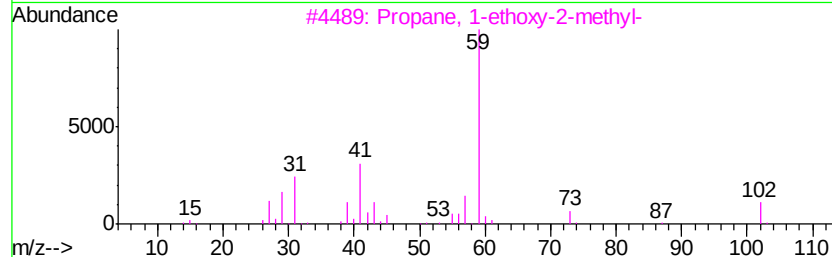
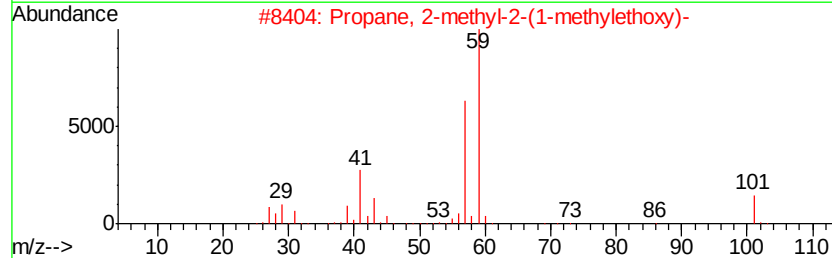
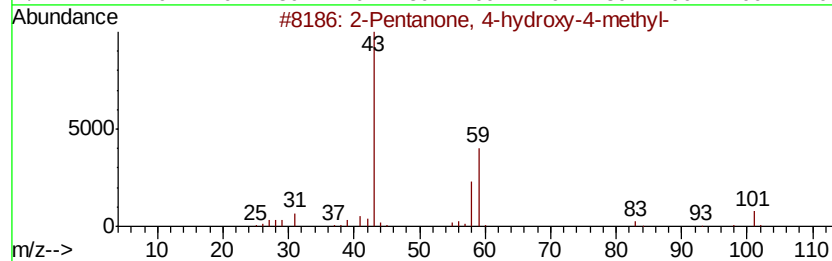
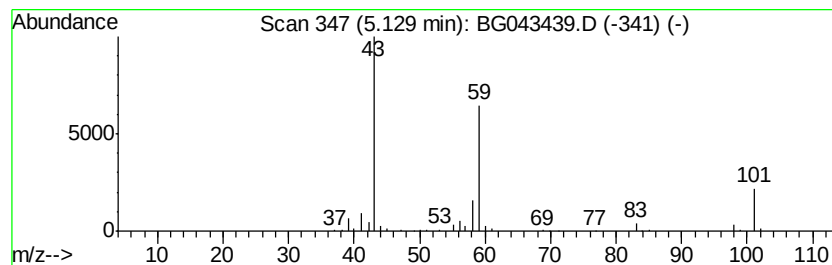
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.13	18.13 ng	224029	1,4-Dichlorobenzene-d4	8.02

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		Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	25
4		1,8-Nonanediol, 8-methyl-	174	C10H22O2	054725-73-4	23
5		Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	17



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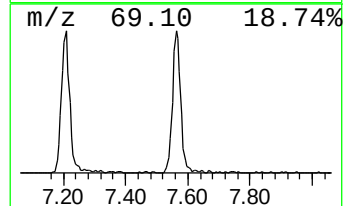
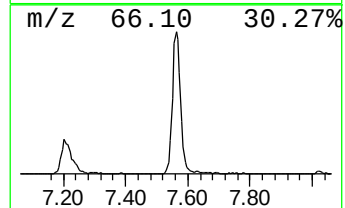
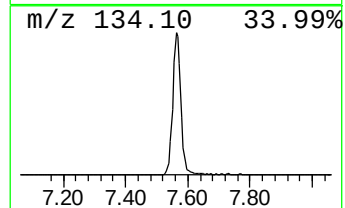
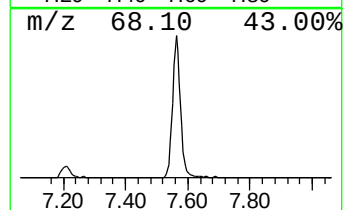
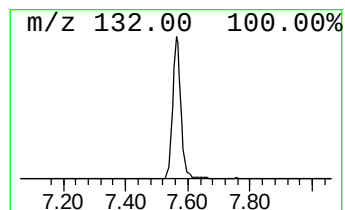
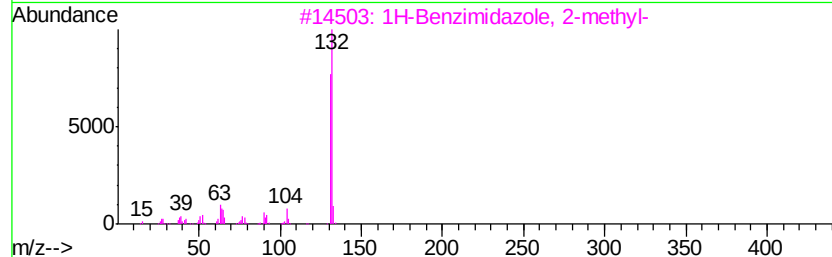
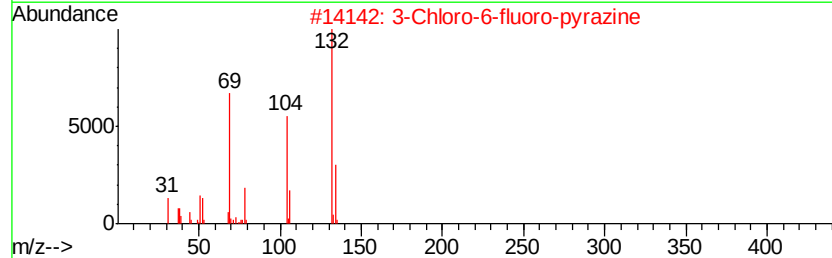
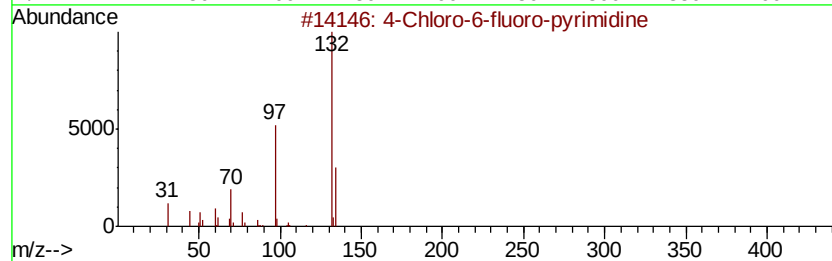
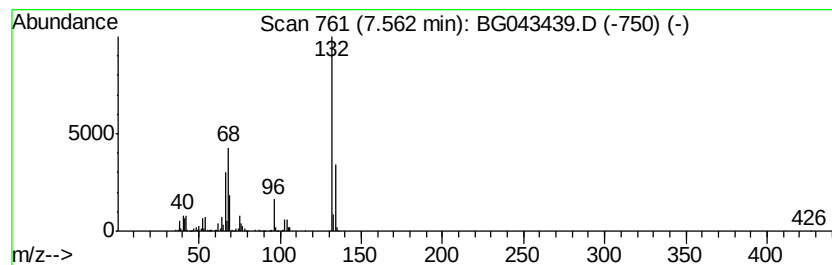
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TIC Library : C:\Database\NIST11.L
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Peak Number 3 unknown7.56 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.56	78.90 ng	974859	1,4-Dichlorobenzene-d4	8.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
3		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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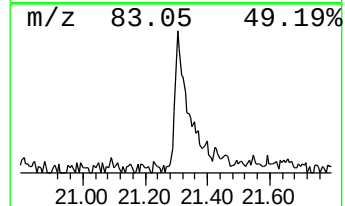
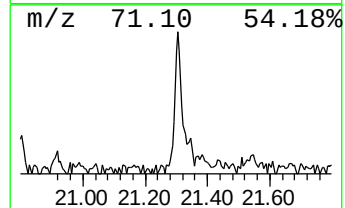
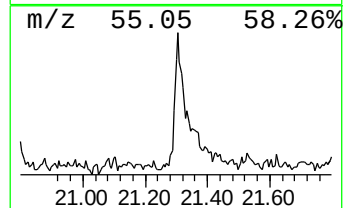
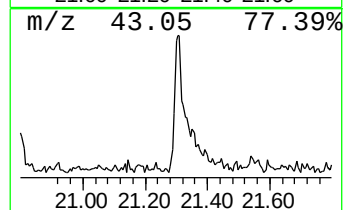
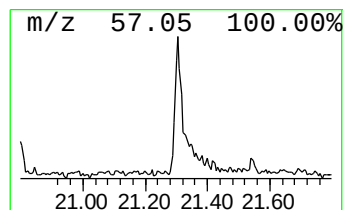
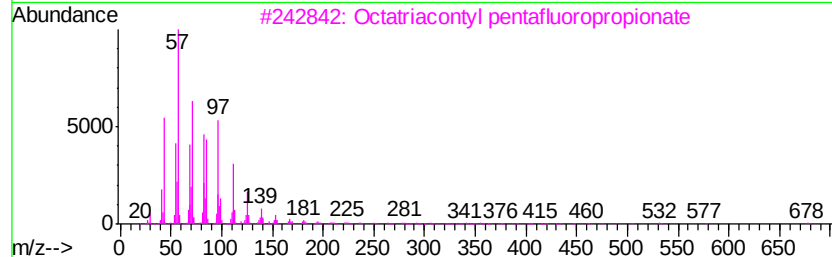
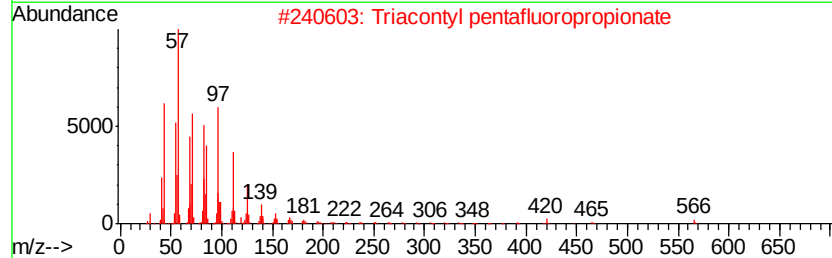
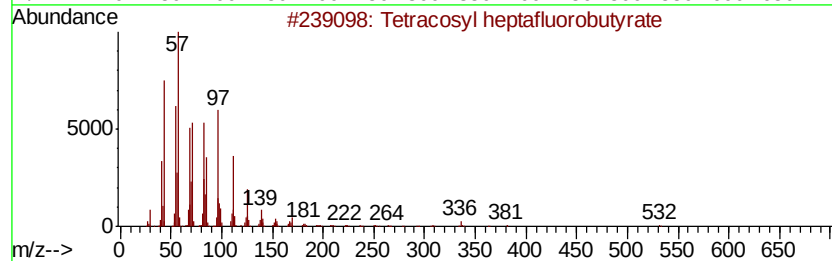
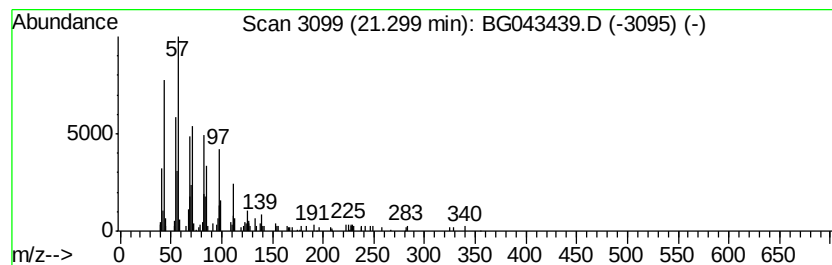
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TIC Library : C:\Database\NIST11.L
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 Peak Number 5 Tetracosyl heptafluorobutyrate Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.30	3.87 ng	177191	Chrysene-d12	21.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetracosyl heptafluorobutyrate	550	C28H49F7O2	1000351-83-7	91
2		Triacetyl pentafluoropropionate	584	C33H61F5O2	1000351-80-0	91
3		Octatriacontyl pentafluoropropio...	697	C41H77F5O2	1000351-89-1	91
4		Heptacosyl pentafluoropropionate	542	C30H55F5O2	1000351-81-6	91
5		Dotriacontyl pentafluoropropionate	612	C35H65F5O2	1000351-81-4	91



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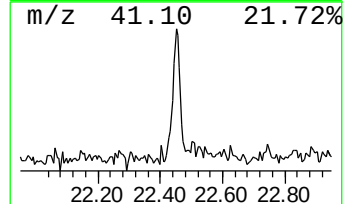
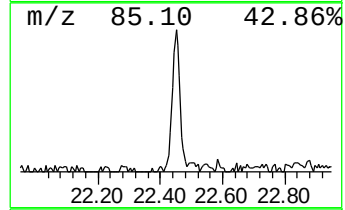
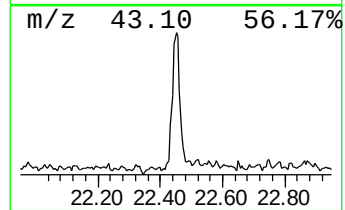
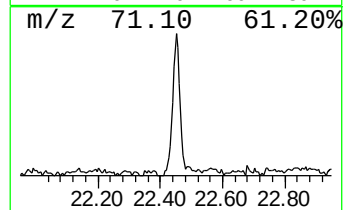
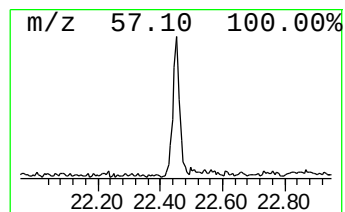
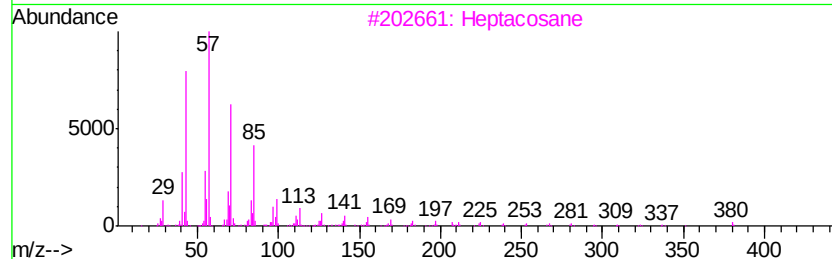
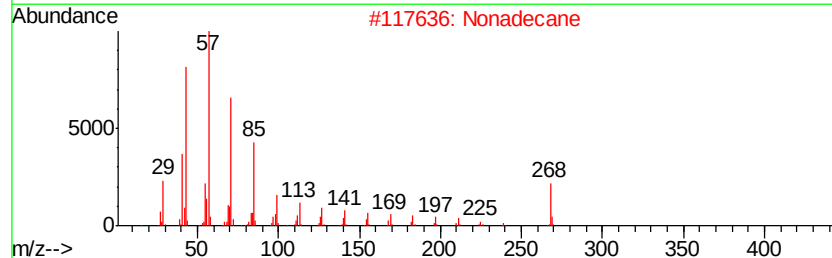
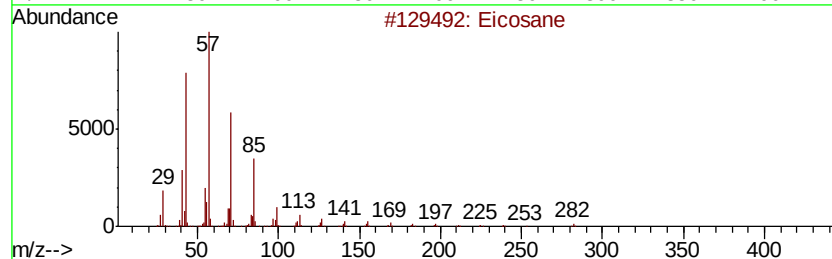
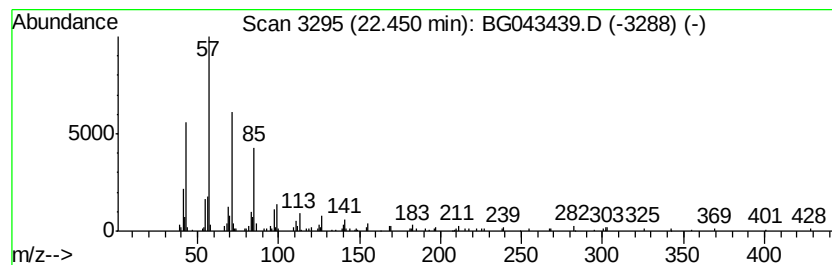
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Peak Number 6 Eicosane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.45	3.06 ng	139959	Chrysene-d12	21.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	95
2		Nonadecane	268	C19H40	000629-92-5	91
3		Heptacosane	380	C27H56	000593-49-7	90
4		Octacosane	394	C28H58	000630-02-4	87
5		Heneicosane	296	C21H44	000629-94-7	87



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Sample : K5587-15
Misc :
ALS Vial : 6 Sample Multiplier: 1

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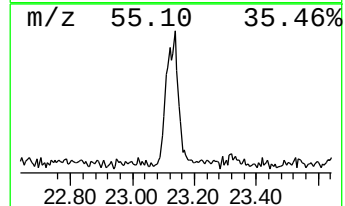
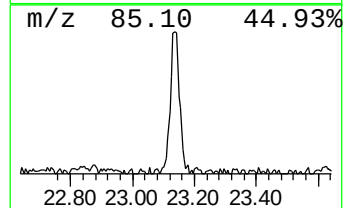
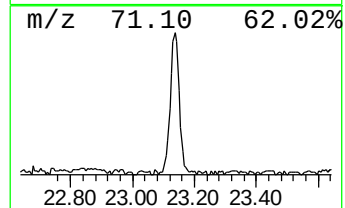
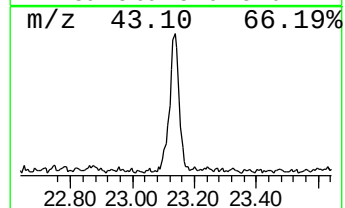
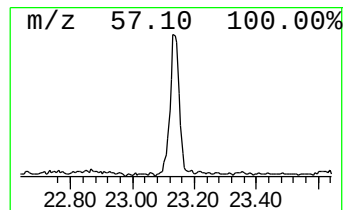
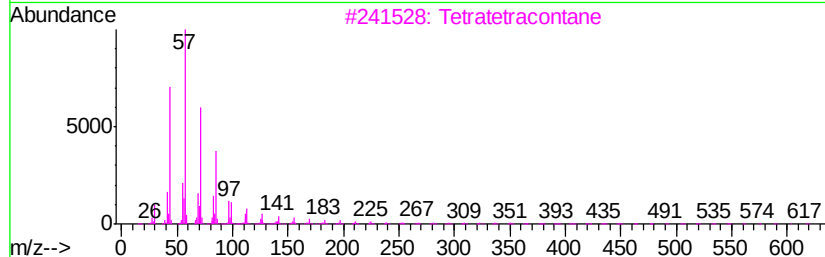
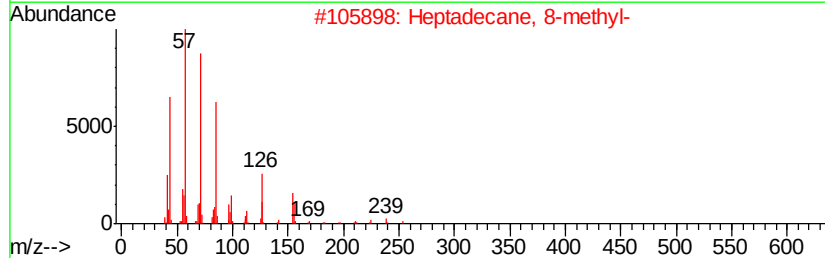
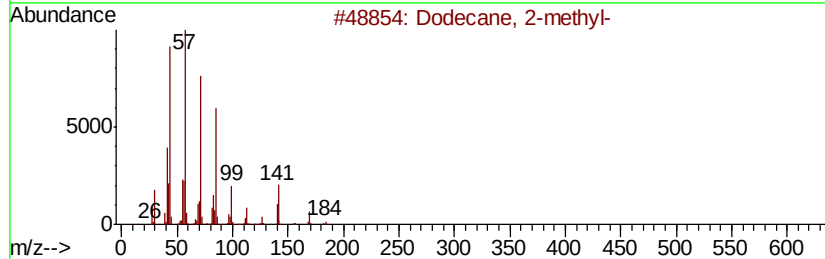
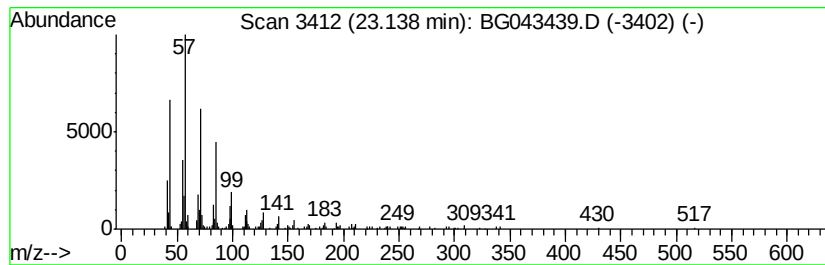
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Dodecane, 2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.14	6.35 ng	290726	Chrysene-d12	21.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Dodecane, 2-methyl-	184	C13H28	001560-97-0	93
2		Heptadecane, 8-methyl-	254	C18H38	013287-23-5	93
3		Tetratetracontane	619	C44H90	007098-22-8	90
4		2-methyloctacosane	408	C29H60	1000376-72-8	90
5		Hentriacontane	437	C31H64	000630-04-6	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG103119\
Data File : BG043439.D
Acq On : 31 Oct 2019 18:24
Operator : JU
Sample : K5587-15
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
CSA-2-1-11

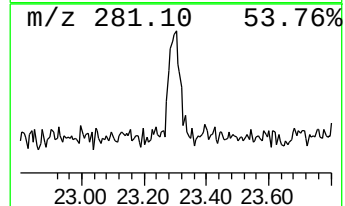
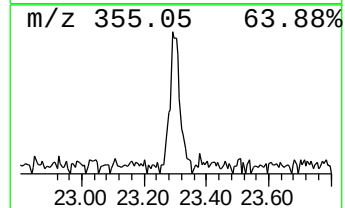
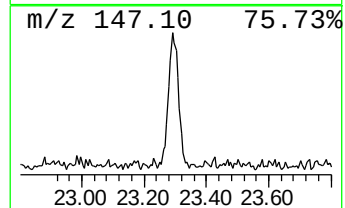
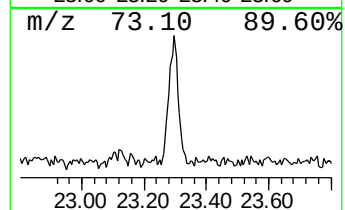
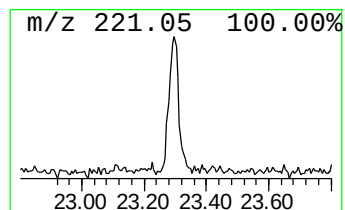
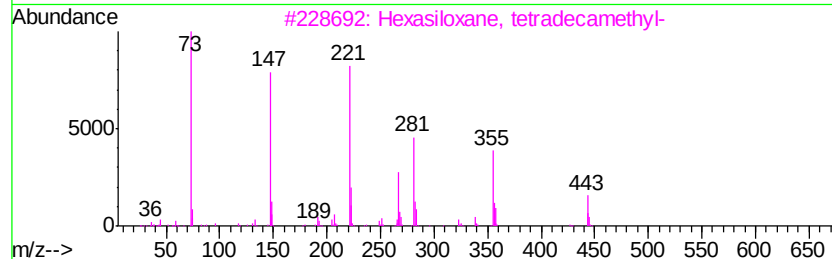
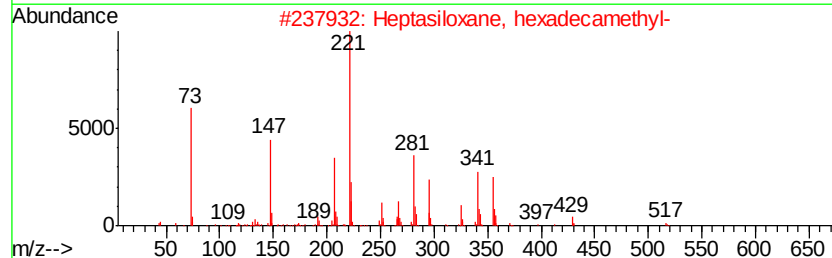
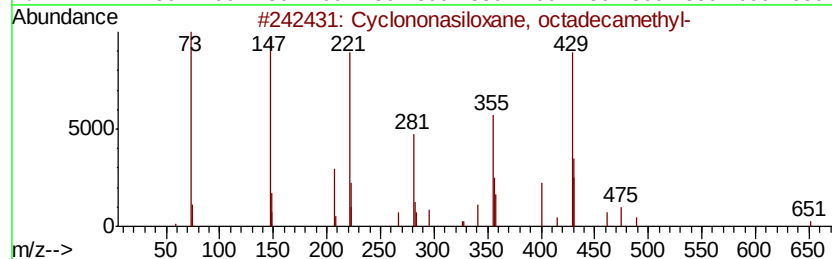
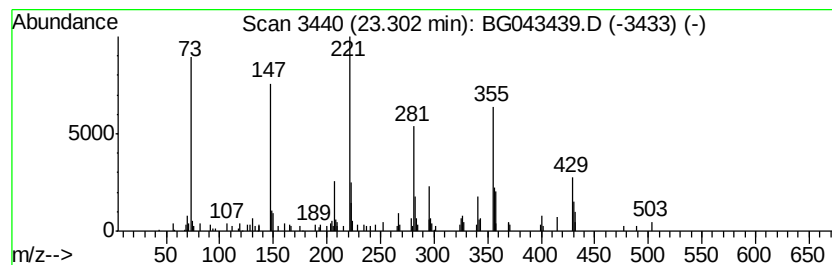
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 unknown23.30 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.30	2.89 ng	160960	Perylene-d12	24.86

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclononasiloxane, octadecamethyl-	666	C18H54O9Si9	000556-71-8	50
2			Heptasiloxane, hexadecamethyl-	532	C16H48O6Si7	000541-01-5	49
3			Hexasiloxane, tetradecamethyl-	458	C14H42O5Si6	000107-52-8	38
4			Cyclodecasiloxane, eicosamethyl-	740	C20H60O10Si10	018772-36-6	38
5			3,6-Dioxa-2,4,5,7-tetrasilaoctan...	294	C10H30O2Si4	004342-25-0	27



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG103119\
 Data File : BG043439.D
 Acq On : 31 Oct 2019 18:24
 Operator : JU
 Sample : K5587-15
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

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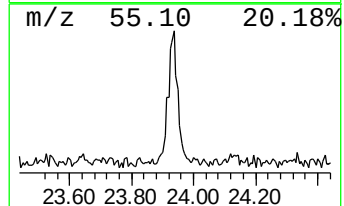
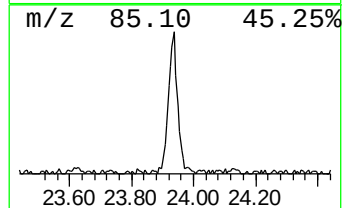
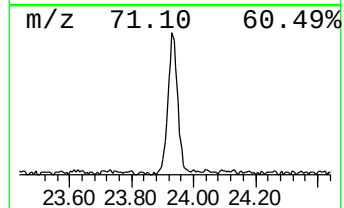
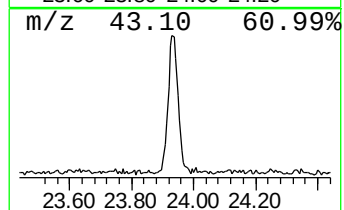
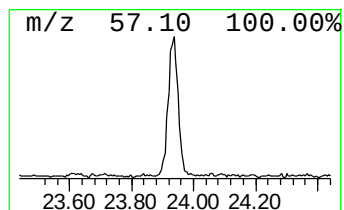
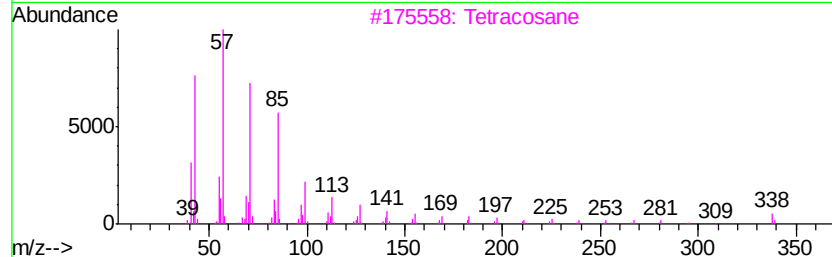
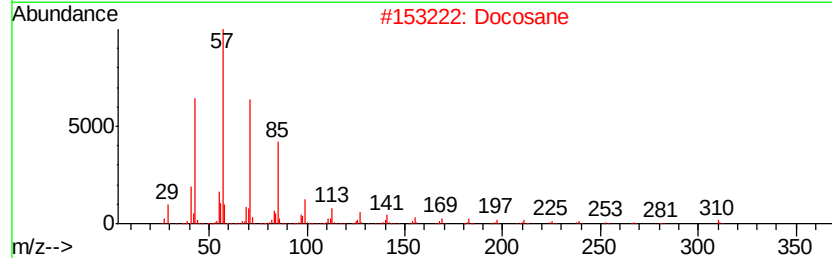
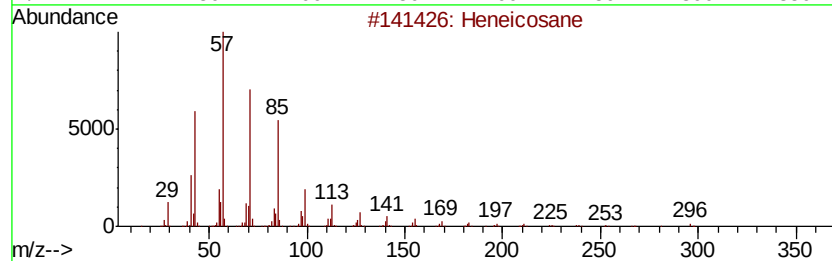
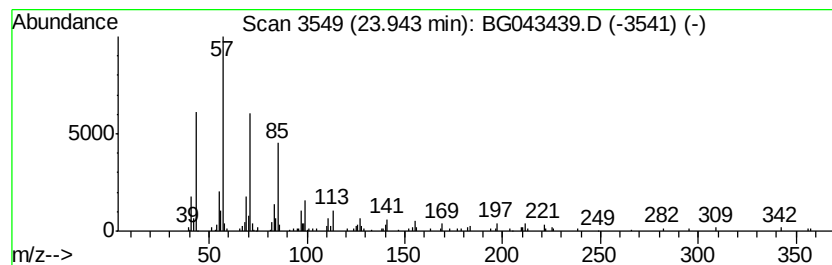
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 9 Heneicosane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.94	4.46 ng	248640	Perylene-d12	24.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	90
2		Docosane	310	C22H46	000629-97-0	87
3		Tetracosane	338	C24H50	000646-31-1	86
4		Eicosane	282	C20H42	000112-95-8	86
5		Octadecane	254	C18H38	000593-45-3	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG103119\
Data File : BG043439.D
Acq On : 31 Oct 2019 18:24
Operator : JU
Sample : K5587-15
Misc :
ALS Vial : 6 Sample Multiplier: 1

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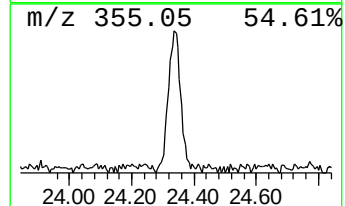
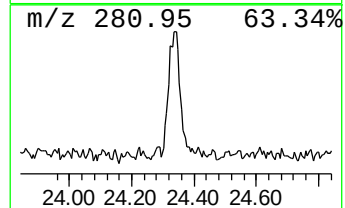
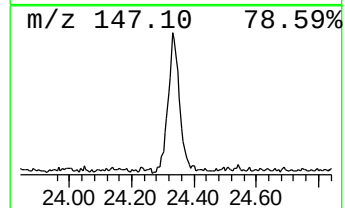
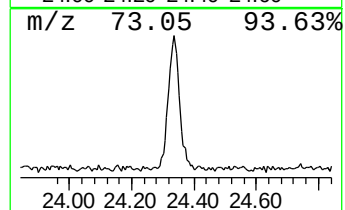
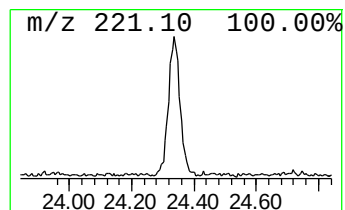
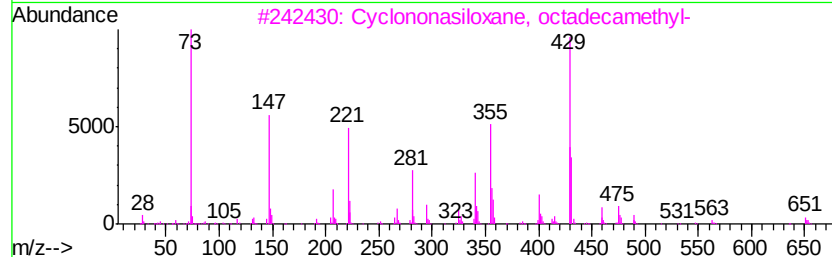
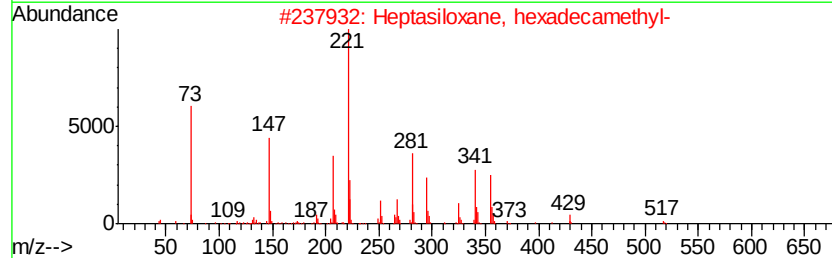
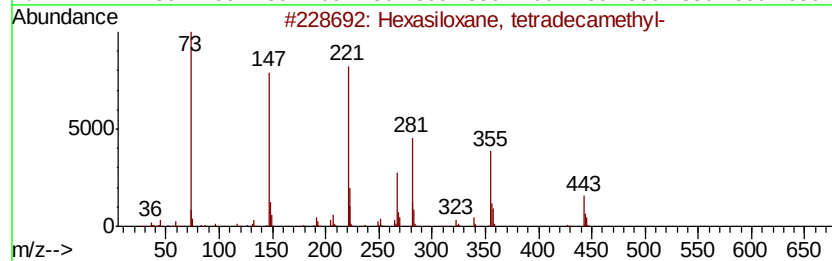
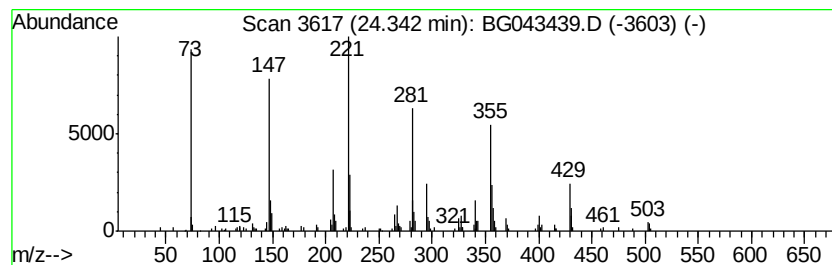
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Hexasiloxane, tetradecamethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.34	5.55 ng	308985	Perylene-d12	24.86

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexasiloxane, tetradecamethyl-	458	C14H42O5Si6	000107-52-8	64
2			Heptasiloxane, hexadecamethyl-	532	C16H48O6Si7	000541-01-5	58
3			Cyclononasiloxane, octadecamethyl-	666	C18H54O9Si9	000556-71-8	25
4			N-Benzyl-N-ethyl-p-isopropylbenz...	281	C19H23NO	015089-22-2	22
5			3,6-Dioxa-2,4,5,7-tetrasilaoctan...	294	C10H30O2Si4	004342-25-0	14



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG103119\
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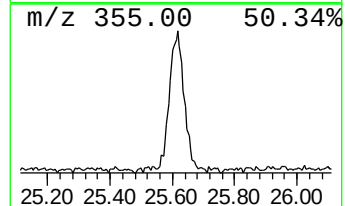
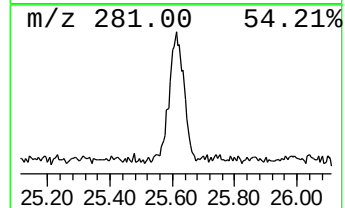
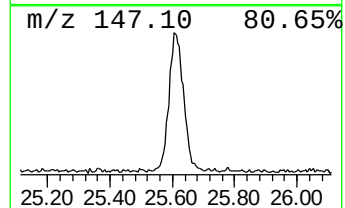
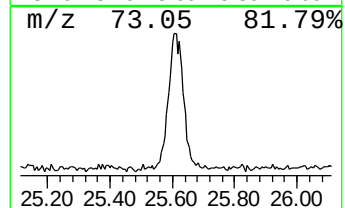
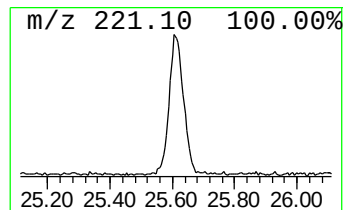
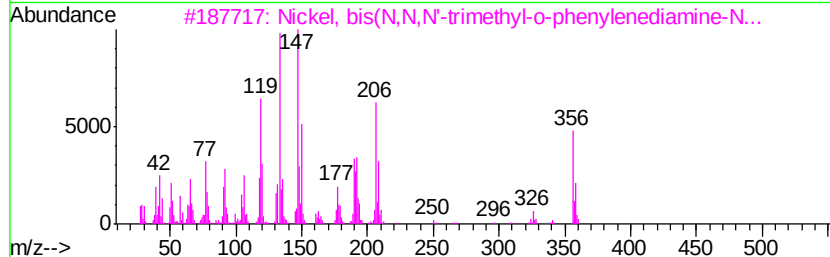
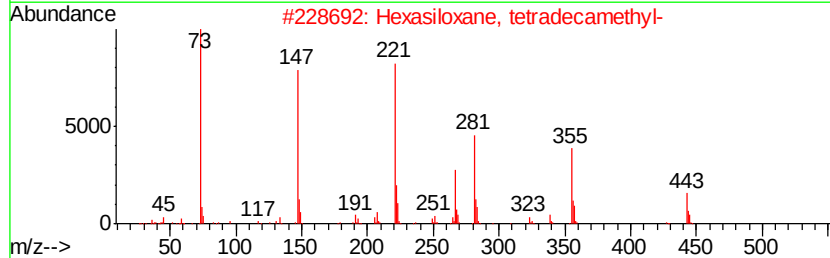
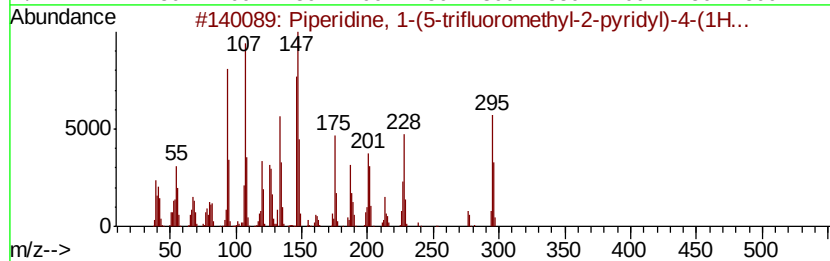
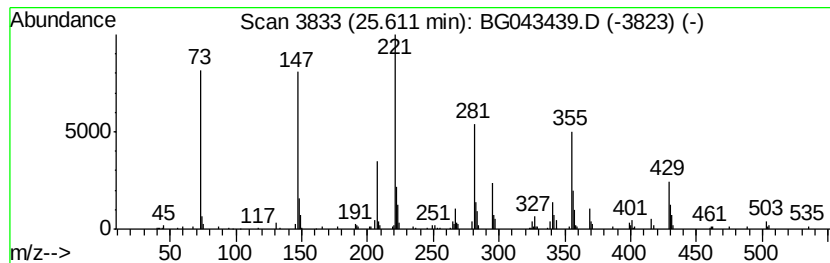
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Piperidine, 1-(5-trifluoromethyl-2-pyridyl)-4-(1H... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
25.61	8.62 ng	480072	Perylene-d12	24.86	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Piperidine, 1-(5-trifluoromethyl-2-pyridyl)-4-(1H-imidazol-5-yl)-	295	C15H16F3N3	1000268-74-7	68
2	Hexasiloxane, tetradecamethyl-	458	C14H42O5Si6	000107-52-8	64
3	Nickel, bis(N,N,N'-trimethyl-o-p-tolyl)-	356	C18H26N4Ni	1000153-51-1	60
4	Cyclononasiloxane, octadecamethyl-	666	C18H54O9Si9	000556-71-8	58
5	Heptasiloxane, hexadecamethyl-	532	C16H48O6Si7	000541-01-5	43



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG103119\
Data File : BG043439.D
Acq On : 31 Oct 2019 18:24
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ClientSampleId :
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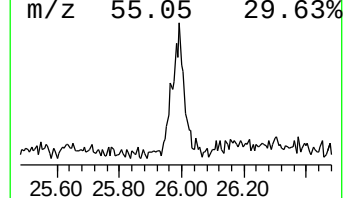
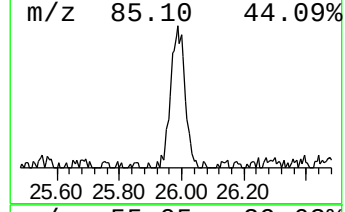
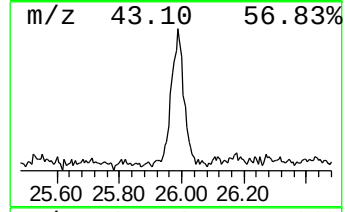
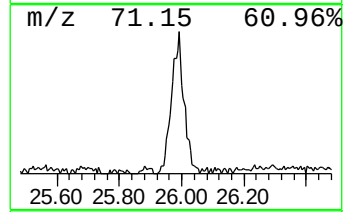
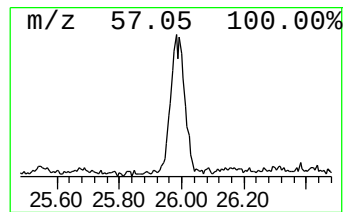
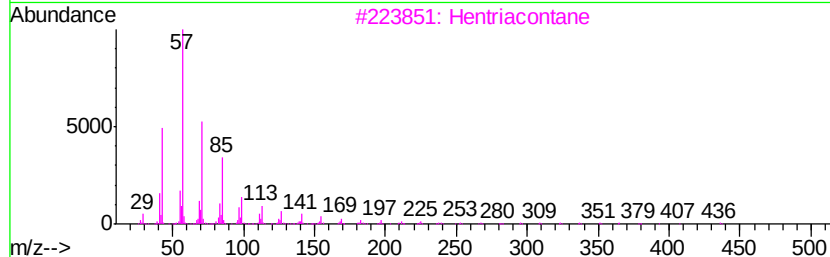
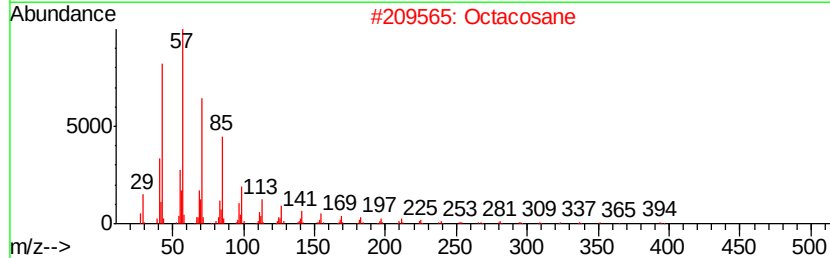
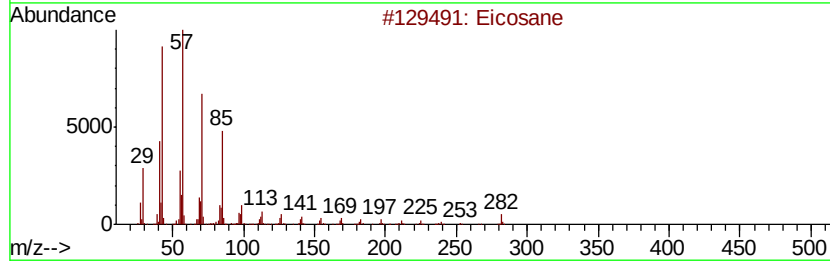
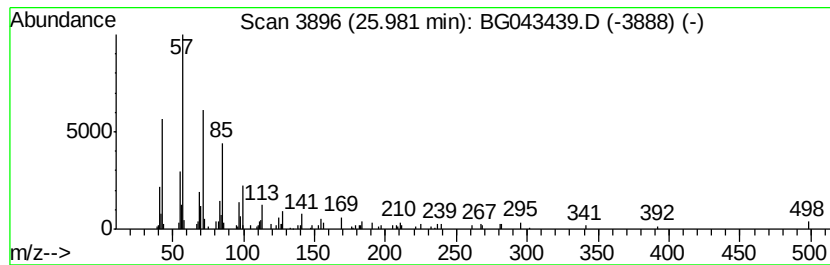
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Octacosane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.98	3.01 ng	167854	Perylene-d12	24.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	95
2		Octacosane	394	C28H58	000630-02-4	87
3		Hentriacontane	437	C31H64	000630-04-6	87
4		Heneicosane	296	C21H44	000629-94-7	87
5		Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	83



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG103119\
Data File : BG043439.D
Acq On : 31 Oct 2019 18:24
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ClientSampleId :
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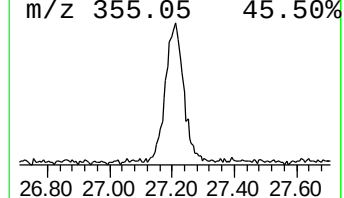
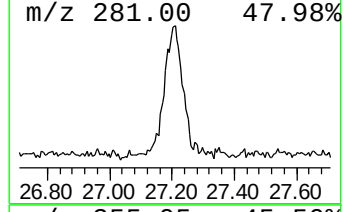
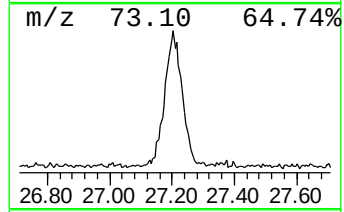
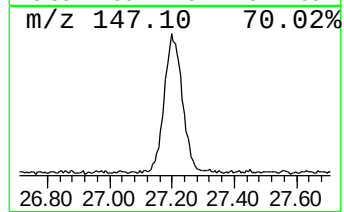
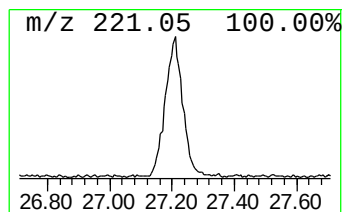
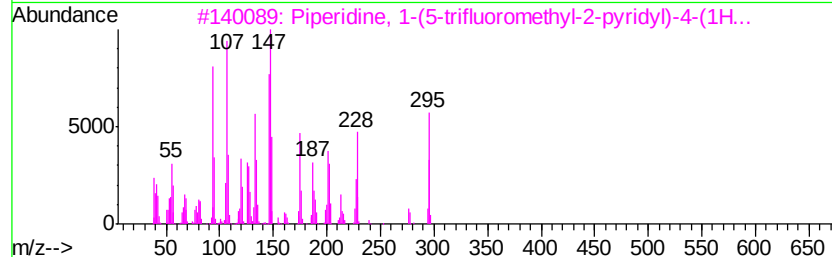
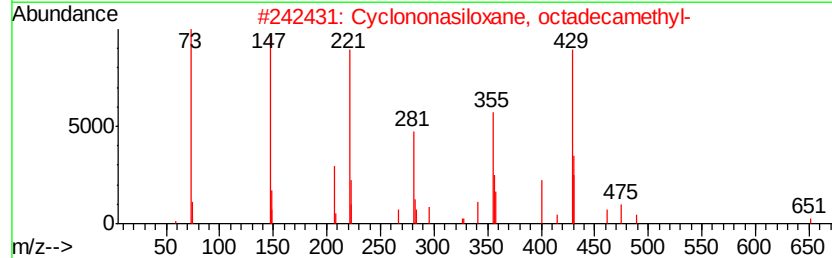
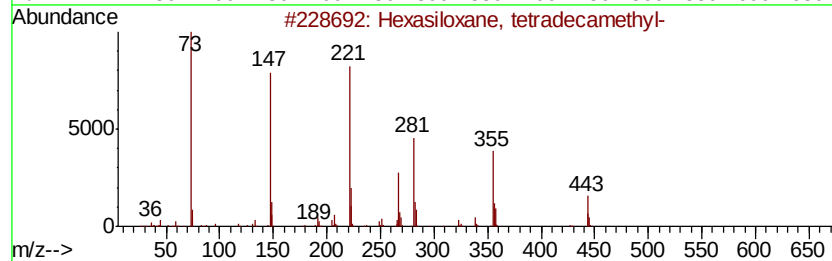
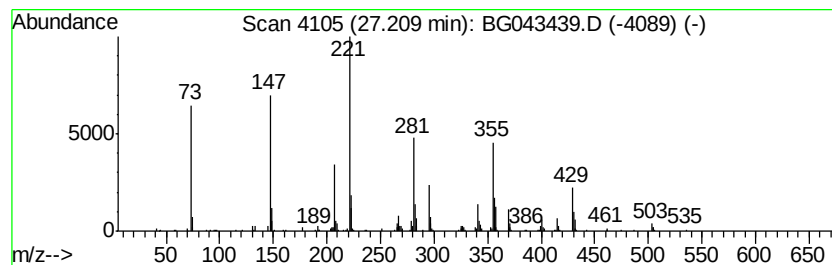
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 unknown27.21 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
27.21	12.63 ng	703632	Perylene-d12	24.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexasiloxane, tetradecamethyl-	458	C14H42O5Si6	000107-52-8	64
2		Cyclononasiloxane, octadecamethyl-	666	C18H54O9Si9	000556-71-8	59
3		Piperidine, 1-(5-trifluoromethyl-2-pyridyl)-4-(1H...	295	C15H16F3N3	1000268-74-7	53
4		2-Amino-2-oxo-acetic acid, N-[3,...	221	C12H15NO3	024451-17-0	35
5		Cyclodecasiloxane, eicosamethyl-	740	C20H60O10Si10	018772-36-6	25



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG103119\
Data File : BG043439.D
Acq On : 31 Oct 2019 18:24
Operator : JU
Sample : K5587-15
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
CSA-2-1-11

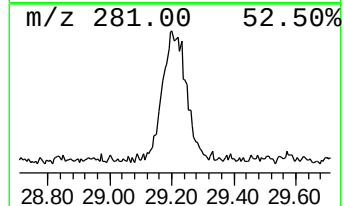
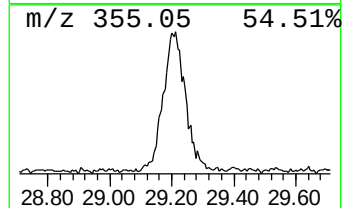
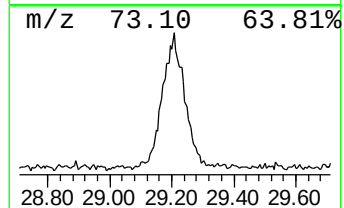
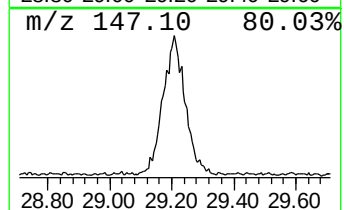
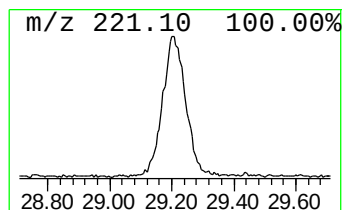
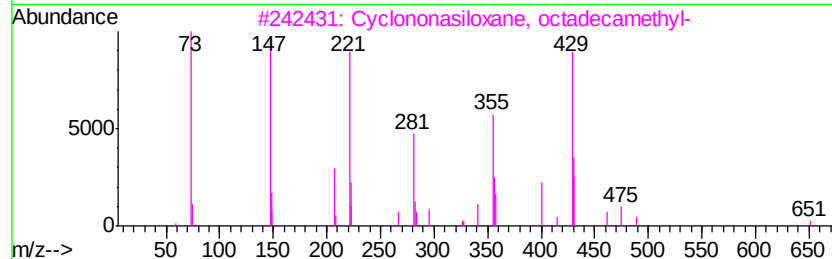
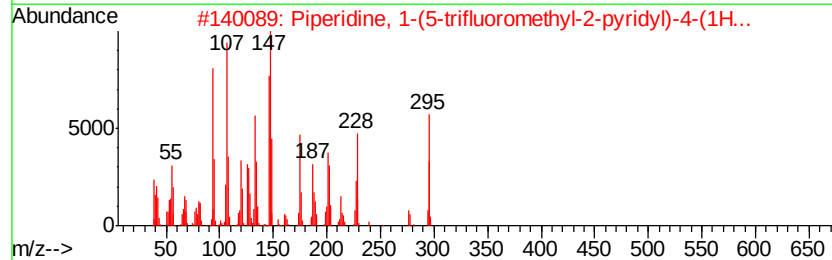
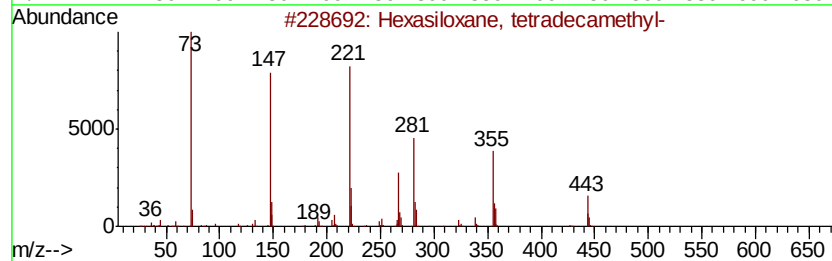
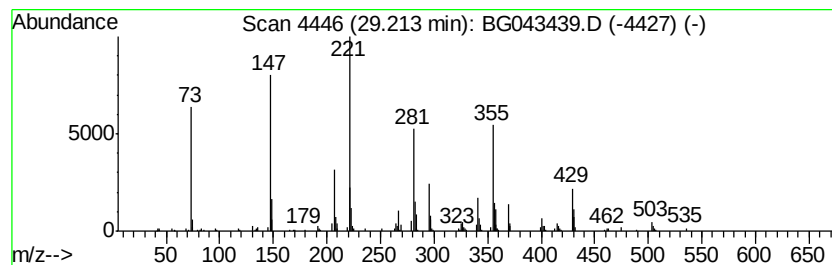
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 unknown29.21 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
29.21	15.24 ng	849099	Perylene-d12	24.86

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexasiloxane, tetradecamethyl-	458	C14H42O5Si6	000107-52-8	64
2			Piperidine, 1-(5-trifluoromethyl-...	295	C15H16F3N3	1000268-74-7	60
3			Cyclononasiloxane, octadecamethyl-	666	C18H54O9Si9	000556-71-8	47
4			Heptasiloxane, hexadecamethyl-	532	C16H48O6Si7	000541-01-5	42
5			N-Benzyl-N-ethyl-p-isopropylbenz...	281	C19H23NO	015089-22-2	30



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG103119\
Data File : BG043439.D
Acq On : 31 Oct 2019 18:24
Operator : JU
Sample : K5587-15
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
CSA-2-1-11

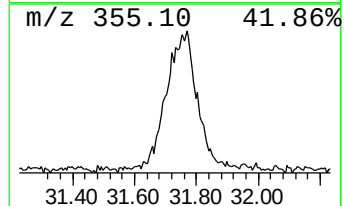
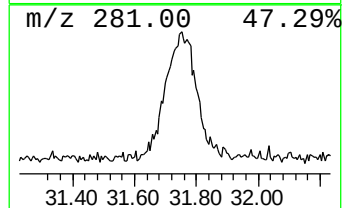
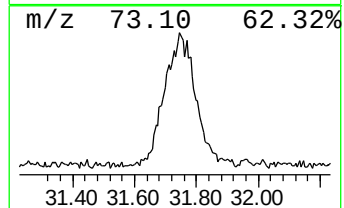
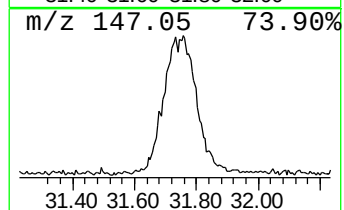
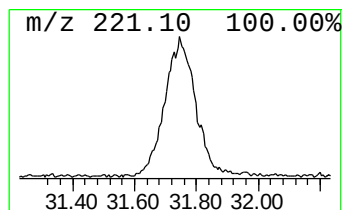
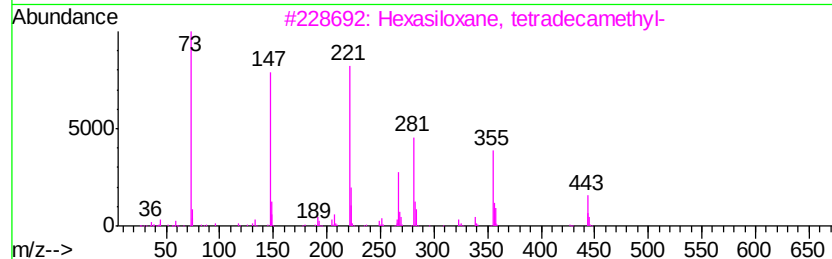
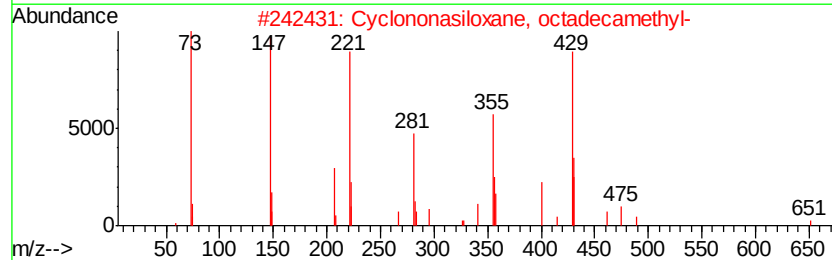
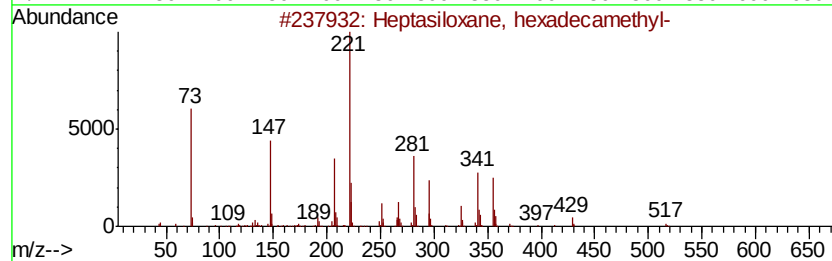
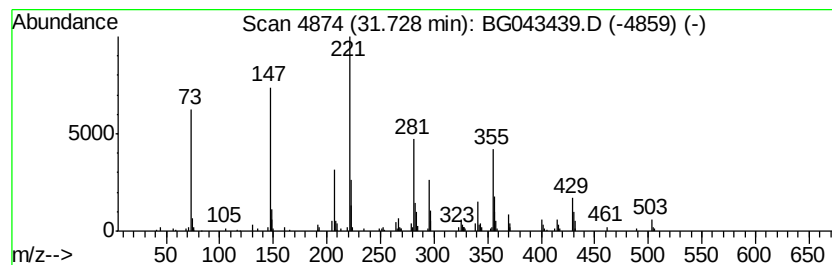
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG102119.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Cyclononasiloxane, octadeca... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
31.73	6.73 ng	375119	Perylene-d12	24.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptasiloxane, hexadecamethyl-	532	C16H48O6Si7	000541-01-5	78
2		Cyclononasiloxane, octadecamethyl-	666	C18H54O9Si9	000556-71-8	74
3		Hexasiloxane, tetradecamethyl-	458	C14H42O5Si6	000107-52-8	72
4		3,6-Dioxa-2,4,5,7-tetrasilaoctan...	294	C10H30O2Si4	004342-25-0	37
5		2-Amino-2-oxo-acetic acid, N-[3,...	221	C12H15N03	024451-17-0	37



Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG103119\
 Data File : BG043439.D
 Acq On : 31 Oct 2019 18:24
 Operator : JU
 Sample : K5587-15
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 CSA-2-1-11

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG102119.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...	3.20	18.5	ng	229127	1	8.02	247114	20.0
2-Pentanone, 4-hy...	5.13	18.1	ng	224029	1	8.02	247114	20.0
unknown7.56	7.56	78.9	ng	974859	1	8.02	247114	20.0
Tetracosyl heptaf...	21.30	3.9	ng	177191	5	21.66	915615	20.0
Eicosane	22.45	3.1	ng	139959	5	21.66	915615	20.0
Dodecane, 2-methyl-	23.14	6.3	ng	290726	5	21.66	915615	20.0
unknown23.30	23.30	2.9	ng	160960	6	24.86	1114460	20.0
Heneicosane	23.94	4.5	ng	248640	6	24.86	1114460	20.0
Hexasiloxane, tet...	24.34	5.5	ng	308985	6	24.86	1114460	20.0
Piperidine, 1-(5-...	25.61	8.6	ng	480072	6	24.86	1114460	20.0
Octacosane	25.98	3.0	ng	167854	6	24.86	1114460	20.0
unknown27.21	27.21	12.6	ng	703632	6	24.86	1114460	20.0
unknown29.21	29.21	15.2	ng	849099	6	24.86	1114460	20.0
Cyclononasiloxane...	31.73	6.7	ng	375119	6	24.86	1114460	20.0