

**SUMMARY OF ANALYSIS (SAMPLE ID: SA38499)**

<b>Testing Location:</b> Arkansas 232 S. Broadview St. Greenbrier, AR 72058 License: ADH 113	<b>Customer ID:</b> 2168 River Valley Relief MIPS 5601 Old Greenwood Rd Suite C Fort Smith, AR 72903 License: 00065C	<b>Order ID:</b> OR11123 <b>Lot Number:</b> M00065C13244752773 <b>Production Run:</b> E20240729BDZLB02	<b>Sample Type:</b> Primary <b>Matrix:</b> Concentrate <b>Mass:</b> 4ea <b>Date Collected:</b> 07/29/2024 <b>Date Received:</b> 07/30/2024
<b>Cultivar (Strain) or Sample Description:</b> Baker's Dozen Hybrid Live Badder 1g Jar			<b>Date Completed:</b> 07/31/2024

\*This page is simply a summary of the analysis performed. For analytical details, please consult the individual Certificate(s) of Analysis for each of the specific test(s) performed. All contaminant action levels are referenced from the State of Arkansas MMJ testing guidelines.

**Moisture Content (%)**

Not Tested

**Water Activity (aw)**

Not Tested

**PASS/FAIL**

**PASS**

Moisture content/water activity action levels are referenced from the State of Arkansas MMJ testing guidelines.

Moisture content levels less than 15% are recommended but the sample does not fail. Water activity levels must be less than 0.65aw.

<b>Cannabinoids (Top 3)</b>	<b>(%)</b>	<b>mg/g</b>
THCA	85.1	851
Δ9-THC	2.37	23.7
CBGA	2.09	20.9
TOTAL CBD	0.0542	0.542
TOTAL THC	77.0	770
TOTAL CANNABINOIDS	90.8	908

<b>Terpenes (Top 5)</b>	<b>(%)</b>	<b>µg/g</b>
β-Caryophyllene	3.03	30300
d-Limonene	3.00	30000
α-Humulene	0.790	7900
Linalool	0.510	5100
β-Myrcene	0.469	4690
TOTAL TERPENES	9.10	91000

<b>Contaminants</b>	<b>PASS/FAIL</b>
Heavy Metals:	<b>PASS</b>
Microbiology:	<b>PASS</b>
Pesticides:	<b>PASS</b>
Residual Solvents:	<b>PASS</b>

**Sample Picture Upon Receipt**



Scan the QR code to verify results.

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*Kyle W. Felling*  
Kyle W. Felling, Ph.D.  
Laboratory Director

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<b>Cultivar (Strain) or Sample Description:</b> Baker's Dozen Hybrid Live Badder 1g Jar			<b>Date Completed:</b> 07/31/2024

**CANNABINOID (POTENCY) PROFILE (SOP: SOP-CANN-001)**

**Analysis Date/Time:** 07/30/2024 1210      **Method:** HPLC/DAD      **Moisture Content (%):** -  
**Analyst:** PW      **Instrument:** Agilent 1100      **Water Activity (aw):** -

<u>Cannabinoid</u>	<u>Result (%)</u>	<u>Result (mg/g)</u>	<u>LOD (mg/g)</u>	<u>LOQ (mg/g)</u>	<u>Result (mg/mL)</u>	<u>Per Serving (mg)</u>	<u>Per Unit (mg)</u>
CBC	ND	ND	0.0944	0.220	-	-	-
CBCA	ND	ND	0.293	0.684	-	-	-
CBD	ND	ND	0.666	1.55	-	-	-
CBDA	0.0618	0.618	0.245	0.572	-	0.618	0.618
CBDV	ND	ND	0.107	0.249	-	-	-
CBDVA	ND	ND	0.285	0.665	-	-	-
CBG	0.178	1.78	0.432	1.01	-	1.78	1.78
CBGA	2.09	20.9	0.613	0.723	-	20.9	20.9
CBL	ND	ND	0.499	1.17	-	-	-
CBN	ND	ND	0.229	0.536	-	-	-
CBNA	ND	ND	0.248	0.577	-	-	-
Δ9-THC	2.37	23.7	0.275	0.641	-	23.7	23.7
Δ8-THC	ND	ND	0.429	1.00	-	-	-
THCA	85.1	851	0.149	0.349	-	851	851
THCV	ND	ND	0.358	0.834	-	-	-
THCVA	1.05	10.5	0.114	0.266	-	10.5	10.5
<b>TOTAL</b>	<b>90.8</b>	<b>908</b>				<b>908</b>	<b>908</b>
<b>TOTAL CBC</b>	-	-				-	-
<b>TOTAL CBD</b>	0.0542	0.542				0.542	0.542
<b>TOTAL CBDV</b>	-	-				-	-
<b>TOTAL CBG</b>	2.01	20.1				20.1	20.1
<b>TOTAL CBN</b>	-	-				-	-
<b>TOTAL THC</b>	77.0	770				770	770
<b>TOTAL THC V</b>	0.911	9.11				9.11	9.11



**SERVING MASS (g):** 1.00  
**SERVINGS/UNIT:** 1

"-" Not detected above LOD.

*Deviations from standard operating procedure:*  
None

*Recoveries for all analyte standards:* 90-110%  
*Replicate Uncertainties:* <5% RSD, <20% RPD  
*Sample/Reagent Blanks:* <RL for all analytes

Values for plant matter are adjusted for moisture content.

Total CBC = (CBCA x 0.877) + CBC  
 Total CBD = (CBDA x 0.877) + CBD  
 Total CBDV = (CBDVA x 0.867) + CBDV  
 Total CBG = (CBGA x 0.878) + CBG  
 Total CBN = (CBNA x 0.876) + CBN  
 Total THC = (THCA x 0.877) + Δ9-THC  
 Total THC V = (THCVA x 0.867) + THC V

Percentage results are reported by mass.

mg/g results are reported as mass component per mass material.

*Abbreviations:* DAD - Diode Array Detector, HPLC - High Pressure Liquid Chromatography, RL - Reporting Limit, RPD - Relative Percent Difference, RSD - Relative Standard Deviation, DET - Detected (less than LOQ), LOD - Limit of Detection, LOQ - Limit of Quantitation, UM - Measurement Uncertainty

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<b>Cultivar (Strain) or Sample Description:</b> Baker's Dozen Hybrid Live Badder 1g Jar			<b>Date Completed:</b> 07/31/2024

**TERPENOID PROFILE**

**Analysis Date/Time:** 07/30/2024 1306  
**Analyst:** KF

**Method:** GC/MS  
**Instrument:** Agilent 7890/5975

**Deviations from SOP:**  
None

<u>Terpene</u>	<u>Result (µg/g)</u>	<u>Result (%)</u>	
α-Bisabolol	850	0.0850	
Camphene	583	0.0583	
δ-3-Carene	ND	-	
β-Caryophyllene	30300	3.03	█
Caryophyllene oxide	215	0.0215	
p-Cymene	ND	-	
Eucalyptol	<LOQ	-	
Geraniol	48.3	0.00483	
Guaiol	ND	-	
α-Humulene	7900	0.790	█
Isopulegol	<LOQ	-	
d-Limonene	30000	3.00	█
Linalool	5100	0.510	█
β-Myrcene	4690	0.469	█
cis-Nerolidol	<LOQ	-	
trans-Nerolidol	4500	0.450	█
α-Ocimene	365	0.0365	
β-Ocimene	166	0.0166	
α-Pinene	1890	0.189	█
β-Pinene	3970	0.397	█
α-Terpinene	<LOQ	-	
γ-Terpinene	<LOQ	-	
Terpinolene	461	0.0461	
<b>TOTAL</b>	<b>91000</b>	<b>9.10</b>	



*Abbreviations:* GC - Gas Chromatography, MS - Mass Spectrometry, RL - Reporting Limit  
*Abbreviations:* ND - Not Detected, LOD - Limit of Detection, LOQ - Limit of Quantitation

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**Reporting Limit (µg/g): 41.4**

"-" Not detected above LOD.

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**Cultivar (Strain) or Sample Description:** Baker's Dozen Hybrid Live Badder 1g Jar **Date Completed:** 07/31/2024

**RESIDUAL SOLVENT PROFILE (SOP: SOP-RS-001)**

**Analysis Date/Time:** 07/30/2024 1440 **Method:** HS/GC/MS **Deviations from SOP:**  
**Analyst:** KF **Instrument:** Agilent 7890/5975 **None**

Solvent	Result (µg/g)	LOD (µg/g)	LOQ (µg/g)	Action Level (µg/g)	Solvent	Result (µg/g)	LOD (µg/g)	LOQ (µg/g)	Action Level (µg/g)
Acetone (67-64-1)	-	28.9	57.7	5000	n-Heptane (142-82-5)	-	28.9	57.7	5000
Acetonitrile (75-5-8)	-	28.9	57.7	410	n-Hexane (110-54-3)	-	10.1	20.2	290
Benzene (71-43-2)	-	0.289	0.577	2	Isobutane (75-28-5)	-	28.9	57.7	5000
n-Butane (106-97-2)	1660	28.9	57.7	5000	Isopropanol (67-63-0)	-	28.9	57.7	5000
1-Butanol (71-36-3)	-	28.9	57.7	5000	Isopropyl acetate (108-21-4)	-	28.9	57.7	5000
2-Butanol (78-92-2)	-	28.9	57.7	5000	Isopropyl benzene (98-82-8)	-	2.89	5.77	70
2-Butanone (78-93-3)	-	28.9	57.7	5000	Methanol (67-56-1)	-	28.9	57.7	3000
Cyclohexane (110-82-7)	-	28.9	57.7	3880	2-Methylbutane (78-78-4)	-	28.9	57.7	5000
1,2-Dimethoxyethane (110-71-4)	-	2.89	5.77	100	Methylene chloride (75-9-2)	-	28.9	57.7	600
N,N-Dimethylacetamide (127-19-5)	-	28.9	57.7	1090	2-Methylpentane (107-83-5)	-	10.1	20.2	290
2,2-Dimethylbutane (75-83-2)	-	10.1	20.2	290	3-Methylpentane (96-10-0)	-	10.1	20.2	290
2,3-Dimethylbutane (79-29-8)	-	10.1	20.2	290	n-Pentane (109-66-0)	-	28.9	57.7	5000
N,N-Dimethylformamide (68-12-2)	-	28.9	57.7	880	1-Pentanol (71-41-0)	-	28.9	57.7	5000
Dimethylsulfoxide (67-68-5)	-	28.9	57.7	5000	n-Propane (74-98-6)	-	28.9	57.7	5000
1,4-Dioxane (123-91-1)	-	28.9	57.7	380	1-Propanol (71-23-8)	-	28.9	57.7	5000
Ethanol (64-17-5)	-	28.9	57.7	5000	Pyridine (110-86-1)	-	10.1	20.2	200
2-Ethoxyethanol (110-80-5)	-	10.1	20.2	160	Tetrahydrofuran (109-99-9)	-	28.9	57.7	720
Ethyl ether (60-29-7)	-	28.9	57.7	5000	Tetramethylene sulfone (126-33-0)	-	10.1	20.2	160
Ethyl acetate (141-78-6)	-	28.9	57.7	5000	Toluene (108-88-3)	-	28.9	57.7	890
Ethyl benzene (100-41-4)	-	28.9	57.7	2170	o-Xylene (95-47-6)	-	28.9	57.7	2170
Ethylene glycol (107-21-1)	-	28.9	57.7	620	m,p-Xylene (108-38-3 or 106-42-3)	-	28.9	57.7	2170
Ethylene oxide (75-21-8)	-	2.89	5.77	50	Xylenes* (1330-20-7)	-	43.3	86.7	2170



**Color Key**

**RESULT < AL** (Green background)  
**RESULT > AL** (Red background)

"DET" detected less than LOQ  
"- " not detected above LOD  
"\*" - o,m,p-Xylene and Ethylbenzene  
Action levels are referenced from the State of Arkansas MMJ testing guidelines.  
A value of "- " for the action level means that analyte is not currently regulated by the regulations referenced above.

Solvent	Synonym(s)	Solvent	Synonym(s)
Acetonitrile	Methyl Cyanide, ACN	Ethylene glycol	1,2-Ethanediol
1-Butanol	n-Butanol, Butyl Alcohol	Isobutane	2-Methylpropane
2-Butanol	sec-Butyl alcohol	Isopropanol	2-Propanol, IPA
2-Butanone	Methyl ethyl ketone, MEK	Isopropyl Acetate	Acetic acid isopropyl ester
1,2-Dimethoxyethane	Monoglyme	Methanol	Methyl alcohol
2,3-Dimethylbutane	Neohexane	2-Methylbutane	Isopentane
2,3-Dimethylbutane	Diisopropyl	Methylene chloride	Dichloromethane
N,N-Dimethylformamide	DMF	2-Methylpentane	Isohexane
Dimethylsulfoxide	DMSO	1-Pentanol	n-Amyl alcohol
2-Ethoxyethanol	Cellosolve, Ethyl glycol	1-Propanol	Propyl alcohol
Ethyl ether	Diethyl ether, Ether	Tetrahydrofuran	THF
Ethyl acetate	EtOAc	Tetramethylene sulfone	Sulfolane
Ethyl benzene	Phenylethane	Xylene	Dimethylbenzene

**Abbreviations:** HS-Headspace, GC-Gas Chromatography, MS-Mass Spectrometry, RL-Reporting Limit, AL-Action Level, CAS-Chemical Abstract Services, LOD - Limit of Detection, LOQ - Limit of Quantification

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**Cultivar (Strain) or Sample Description:** Baker's Dozen Hybrid Live Badder 1g Jar **Date Completed:** 07/31/2024

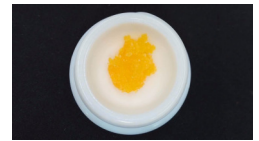
**PESTICIDES PROFILE (SOP: SOP-PEST-001)**

**Analysis Date/Time:** 07/30/2024 1200  
**Analyst:** KF

**Method:** LC/MS/MS  
**Instrument:** Shimadzu LC-8050

**Deviations from SOP:**  
None

Pesticide	Result (µg/g)	LOD (µg/g)	LOQ (µg/g)	Action Level (µg/g)	Pesticide	Result (µg/g)	LOD (µg/g)	LOQ (µg/g)	Action Level (µg/g)
Abamectin (71751-41-2)	-	0.0408	0.327	0.5	Kresoxim-methyl (143390-89-0)	-	0.0408	0.327	0.4
Acephate (30560-19-1)	-	0.0408	0.327	0.4	Malathion (121-75-5)	-	0.0408	0.327	0.2
Acetamiprid (135410-20-7)	-	0.0408	0.327	0.2	Metalaxyl (57837-19-1)	-	0.0408	0.327	0.2
Aldicarb (116-06-3)	-	0.0408	0.327	0.4	Methiocarb (2032-65-7)	-	0.0408	0.327	0.2
Azoxystrobin (131860-33-8)	-	0.0408	0.327	0.2	Methomyl (16752-77-5)	-	0.0408	0.327	0.4
Bifenazate (149877-41-8)	-	0.0408	0.327	0.2	Methyl parathion (298-0-0)	-	0.0408	0.327	0.2
Bifenthrin (82657-04-3)	-	0.0408	0.327	0.2	MGK 264 (113-48-4)	-	0.0408	0.327	0.2
Boscalid (188425-85-6)	-	0.0408	0.327	0.4	Myclobutanil (88671-89-0)	-	0.0408	0.327	0.2
Carbaryl (63-25-2)	-	0.0408	0.327	0.2	Naled (300-76-5)	-	0.0408	0.327	0.5
Carbofuran (1563-66-2)	-	0.0408	0.327	0.2	Oxamyl (23135-22-0)	-	0.0408	0.327	1
Chlorantraniliprole (800008-45-7)	-	0.0408	0.327	0.2	Pacllobutrazol (76738-62-0)	-	0.0408	0.327	0.4
Chlorfenapyr (122453-73-0)	-	0.0408	0.327	1	Permethrins (52645-53-1)	-	0.0408	0.327	0.2
Chlorpyrifos (2921-88-2)	-	0.0408	0.327	0.2	Phosmet (732-11-6)	-	0.0408	0.327	0.2
Clofentezine (74115-24-5)	-	0.0408	0.327	0.2	Piperonyl butoxide (51-03-6)	-	0.0408	0.327	2
Cyfluthrin (68359-37-5)	-	0.0408	0.327	1	Prallethrins (2331-36-9)	-	0.0408	0.327	0.2
Cypermethrin (52315-07-8)	-	0.0408	0.327	1	Propiconazole (60207-90-1)	-	0.0408	0.327	0.4
Daminozide (1596-84-5)	-	0.0408	0.327	1	Propoxur (114-26-1)	-	0.0408	0.327	0.2
DDVP (62-73-7)	-	0.0408	0.327	0.1	Pyrethrins (8003-34-7)	-	0.0408	0.327	1
Diazinon (333-41-5)	-	0.0408	0.327	0.2	Pyridaben (96489-71-3)	-	0.0408	0.327	0.2
Dimethoate (60-51-5)	-	0.0408	0.327	0.2	Spinosad (168316-95-8)	-	0.0408	0.327	0.2
Ethoprophos (13194-48-4)	-	0.0408	0.327	0.2	Spiromesifen (283594-90-1)	-	0.0408	0.327	0.2
Etofenprox (80844-07-1)	-	0.0408	0.327	0.4	Spirotetramat (203313-25-1)	-	0.0408	0.327	0.2
Etoxazole (153233-91-1)	-	0.0408	0.327	0.2	Spiroxamine (118134-30-8)	-	0.0408	0.327	0.4
Fenoxycarb (72490-01-8)	-	0.0408	0.327	0.2	Tebuconazole (80443-41-0)	-	0.0408	0.327	0.4
(E)-Fenpyroximate (134098-61-6)	-	0.0408	0.327	0.4	Thiacloprid (111988-49-9)	-	0.0408	0.327	0.2
Fipronil (120068-37-3)	-	0.0408	0.327	0.4	Thiamethoxam (153719-23-4)	-	0.0408	0.327	0.2
Fonicamid (158062-67-0)	-	0.0408	0.327	1	Trifloxystrobin (141517-21-7)	-	0.0408	0.327	0.2
Fludioxinil (131341-86-1)	-	0.0408	0.327	0.4					
Hexythiazox (78587-05-0)	-	0.0408	0.327	1					
Imazalil (35554-44-0)	-	0.0408	0.327	0.2					
Imidacloprid (138261-41-3)	-	0.0408	0.327	0.4					



**Color Key**

**RESULT < AL**  
**RESULT > AL**

"DET" detected less than LOQ  
"-." not detected above LOD  
Permethrins measured as the cumulative residue of the *cis*- and *trans*- permethrin isomers.  
Pyrethrins measured as the cumulative residue of the pyrethrin I, cinerin I, and jasmolin I isomers.  
Action levels are referenced from the State of Arkansas MMJ testing guidelines.  
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Abbreviations: LC - Liquid Chromatography, MS - Mass Spectrometry, RL - Reporting Limit, AL - Action Level, CAS - Chemical Abstract Services, LOD - Limit of Detection, LOQ - Limit of Quantification

Pesticide	Synonym(s)	Pesticide	Synonym(s)	Pesticide	Synonym(s)
Cyfluthrin	Baythroid	Myclobutanil	Systhane	Propiconazole	Tilt
DDVP	Dichlorvos	Naled	Dibrom	Propoxur	Baygon
Ethoprophos	Prophos	Phosmet	Imidan		

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Laboratory Director



**CERTIFICATE OF ANALYSIS (SAMPLE ID: SA38499)**

<b>Testing Location:</b> Arkansas 232 S. Broadview St. Greenbrier, AR 72058 License: ADH 113	<b>Customer ID:</b> 2168 River Valley Relief MIPS 5601 Old Greenwood Rd Suite C Fort Smith, AR 72903 License: 00065C	<b>Order ID:</b> OR11123 <b>Lot Number:</b> M00065C13244752773 <b>Production Run:</b> E20240729BDZLB02	<b>Sample Type:</b> Primary <b>Matrix:</b> Concentrate <b>Mass:</b> 4ea <b>Date Collected:</b> 07/29/2024 <b>Date Received:</b> 07/30/2024
<b>Cultivar (Strain) or Sample Description:</b> Baker's Dozen Hybrid Live Badder 1g Jar			<b>Date Completed:</b> 07/31/2024

**HEAVY METAL PROFILE (SOP: SOP-ICP-200.6)**

**Analysis Date/Time:** 07/30/2024 1336      **Method:** ICP/MS      **Deviations from SOP:** None  
**Analyst:** KF      **Instrument:** Agilent 7500ce

<u>Heavy Metal</u>	<u>Result (µg/kg)</u>	<u>LOD (µg/kg)</u>	<u>LOQ (µg/kg)</u>	<u>Action Level (µg/kg)</u>
Arsenic (As)	-	59.4	94.1	200
Cadmium (Cd)	-	59.4	94.1	200
Lead (Pb)	-	59.4	94.1	500
Mercury (Hg)	-	59.4	94.1	100



*Abbreviations:* ICP - Inductively-Coupled Plasma, OES - Optical Emission Spectroscopy, DMA - Direct Mercury Analyzer, RL - Reporting Limit, AL - Action Level, LOD - Limit of Detection, LOQ - Limit of Quantitation

**Color Key**



"DET" detected less than LOQ  
 "-" not detected above LOD

Action levels for heavy metals are referenced from the State of Arkansas MMJ testing guidelines.

*Disclaimer:* This information is provided as a service and makes no claims of efficacy and/or safety of this product. Results are applicable only for the sample(s) analyzed and for the specific analysis conducted. This report is for informational purposes only and should not be used to diagnose, treat, or prevent any medical-related symptoms. The statements and results herein have not been approved and/or endorsed by the FDA.

*Kyle W. Felling*  
 Kyle W. Felling, Ph.D.  
 Laboratory Director

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**CERTIFICATE OF ANALYSIS (SAMPLE ID: SA38499)**

<b>Testing Location:</b> Arkansas 232 S. Broadview St. Greenbrier, AR 72058 License: ADA 05_H273	<b>Customer ID:</b> 2168 River Valley Relief MIPS 5601 Old Greenwood Rd Suite C Fort Smith, AR 72903 License: 00065C	<b>Sample ID:</b> SA38499 <b>Lot Number:</b> M00065C13244752773 <b>Production Run:</b> E20240729BDZLB02	<b>Sample Type:</b> Primary <b>Matrix:</b> Concentrate <b>Mass:</b> 4ea <b>Date Collected:</b> 07/29/2024 <b>Date Received:</b> 07/30/2024
<b>Cultivar (Strain) or Sample Description:</b> Baker's Dozen Hybrid Live Badder 1g Jar		<b>Date Completed:</b> 07/31/2024	

**MICROBIOLOGICAL PROFILE (SOP: SOP-Micro-001)**

**Analysis Date/Time:** 07/31/2024 0754      **Method:** Hardy Diagnostics CompactDry      **Deviations from SOP:**  
**Analyst:** PW      **Instrument:** Thermo Incubator      None

<b>Bacteria/Microbe</b>	<b>Result (CFU/g)</b>	<b>Action Level (CFU/g)</b>
Aerobic Plate Count	NT	-
Coliforms, Total	Absent	1
Escherichia Coli (E. Coli)	Absent	100
Mold/Yeast	NT	-
Pseudomonas aeruginosa	NT	-
Salmonella spp.	NT	-
Staphylococcus aureus	NT	-



**Abbreviations:** EC - Escherichia Coli, CFU - Colony-Forming Unit, RL - Reporting Limit, AL - Action Level, TNTC - Too Numerous To Count, NT - Not Tested  
Absent - Not Detected Above RL, Present - Detected Above RL

**Color Key**

**RESULT < AL**

**RESULT > AL**

**Reporting Limit (CFU/g)**  
1

Action levels for microbiology are referenced from the State of Arkansas MMJ testing guidelines.  
A value of "-" for the action level means that analyte is not currently regulated by the regulations referenced above.

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*Kyle W. Felling*  
Kyle W. Felling, Ph.D.  
Laboratory Director

