

Testing Accreditation #: 77802

Test Certificate #: 128457-001

Client Name, Sample Details
Vision Aura Hemp Oil

Sample: Full Spectrum CBD
Type: Concentrate
Method: FE-52 (EN 15662 & AOAC 2007.01)

Test Conditions
Prepsheet ID#: MIPS201226b
Scale: XS205-M12
Temp: 21.2 °C
Baro PE: 972.2 hPa
Analyst: MEH
Technician: ANJ

Sample ID#: 128457
Harvest/Process Date: 12/23/2020
Serving Size (g): 1
Date Received: 12/23/2020
Test Date: 12/26/2020
Valid Through: 12/26/2021
Report Issued: 12/29/2020



Compound	MRL (µg/g)	LOD (µg/g)	Status (µg/g)	Compound	MRL (µg/g)	LOD (µg/g)	Status (µg/g)
Aldicarb	0.400	0.125	Pass/<LOD	Abamectin****	0.500	0.125	Pass/<LOD
Acephate	0.400	0.125	Pass/<LOD	Acequinocyl	2.000	0.250	Pass/<LOD
Acetamiprid	0.200	0.125	Pass/<LOD	Azoxystrobin	0.200	0.125	Pass/<LOD
Bifenazate	0.200	0.125	Pass/<LOD	Bifenthrin	0.200	0.125	Pass/<LOD
Boscalid	0.400	0.125	Pass/<LOD	Carbaryl	0.200	0.125	Pass/<LOD
Carbofuran	0.200	0.125	Pass/<LOD	Chlorantraniliprole	0.200	0.125	Pass/<LOD
Chlorfenapyr	1.000	0.500	Pass/<LOD	Chlorpyrifos	0.200	0.125	Pass/<LOD
Clofentezine	0.200	0.125	Pass/<LOD	Cyfluthrin**	1.000	0.500	Pass/<LOD
Cypermethrin***	1.000	0.500	Pass/<LOD	Daminozide	1.000	0.500	Pass/<LOD
DDVP (Dichlorvos)	1.000	0.250	Pass/<LOD	Diazinon	0.200	0.125	Pass/<LOD
Dimethoate	0.200	0.125	Pass/<LOD	Ethoprophos	0.200	0.125	Pass/<LOD
Etofenprox	0.400	0.125	Pass/<LOD	Etoxazole	0.200	0.125	Pass/<LOD
Fenoxycarb	0.200	0.125	Pass/<LOD	Fenpyroximate	0.400	0.125	Pass/<LOD
Fipronil	0.400	0.125	Pass/<LOD	Flonicamid	1.000	0.125	Pass/<LOD
Fludioxonil	0.400	0.125	Pass/<LOD	Hexythiazox	1.000	0.125	Pass/<LOD
Imazalil	0.200	0.125	Pass/<LOD	Imidacloprid	0.400	0.125	Pass/<LOD
Kresoxim Methyl	0.400	0.125	Pass/<LOD	Malathion	0.200	0.125	Pass/<LOD
Metalaxyl	0.200	0.125	Pass/<LOD	Methiocarb	0.200	0.125	Pass/<LOD
Methomyl	0.400	0.125	Pass/<LOD	Methyl Parathion	0.200	0.125	Pass/<LOD
MGK-264‡	0.200	0.125	Pass/<LOD	Myclobutanil	0.200	0.125	Pass/<LOD
Naled	0.500	0.125	Pass/<LOD	Oxamyl	1.000	0.125	Pass/<LOD
Paclobutrazol	0.400	0.125	Pass/<LOD	Permethrins†	0.200	0.125	Pass/<LOD
Phosmet	0.200	0.125	Pass/<LOD	Piperonyl Butoxide	100.000	1.900	Pass/<LOD
Prallethrin	0.200	0.125	Pass/<LOD	Propiconazole	0.400	0.125	Pass/<LOD
Propoxur	0.200	0.125	Pass/<LOD	Pyrethrins*	1.000	0.125	Pass/<LOD
Pyridaben	0.200	0.125	Pass/<LOD	Spinosad*****	0.200	0.125	Pass/<LOD
Spiromesifen	0.200	0.125	Pass/<LOD	Spirotetramat	0.200	0.125	Pass/<LOD
Spiroxamine‡	0.400	0.125	Pass/<LOD	Tebuconazole	0.400	0.125	Pass/<LOD
Thiacloprid	0.200	0.125	Pass/<LOD	Thiamethoxam	0.200	0.125	Pass/<LOD
Trifloxystrobin	0.200	0.125	Pass/<LOD				

* Pyrethrins are reported as the sum of Jasmolin I, Cinerin I, and Pyrethrin I

** Cyfluthrins are reported as the sum of isomers Cyfluthrin I, II, III, and IV

*** Cypermethrins are reported as the sum of isomers Cypermethrin I, II, III, and IV

**** Abamectin is reported as the sum of Avermectin B1a and Avermectin B1b

***** Spinosad is reported as the sum of Spinosyn A and Spinosyn D

† Permethrin and Prallethrin are reported as the sum of cis and trans isomers

‡ MGK-264 and Spiroximine are reported as the sum of isomers I and II

MRL - Maximum Residue Limit; LOD - Limit of Detection

Sample was sampled and tested in accordance with the Safety Compliance Facility Sampling and Testing Information.

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Katrina Barnes, Lab Manager




Mackenzie E. Hyman, Quality Manager

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Tested by Iron Laboratories Michigan, 1825 E. West Maple Walled Lake, MI 48390

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Vision Aura Hemp Oil

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Target Compound Name	Method Blank (µg/g)	QC Spike (µg/g)	Matrix Spike (µg/g)	Matrix Spike Duplicate (µg/g)	MS recovery%	MSD recovery%	Relative Percent Difference (%)	QC Flag
Acephate	N.D.	1	0.971	0.992	97.10	99.20	2.14	
Acequinocyl	N.D.	1	0.26	0.288	26.00	28.80	10.22	LR
Acetamiprid	N.D.	1	0.888	0.916	88.80	91.60	3.10	
Aldicarb	N.D.	1	0.961	0.944	96.10	94.40	1.78	
Avermectin B1a	N.D.	0.97	1.49	0.992	153.61	102.27	40.13	HRQ
Azoxystrobin	N.D.	1	0.958	0.961	95.80	96.10	0.31	
Bifenazate	N.D.	1	0.447	0.446	44.70	44.60	0.22	LR
Bifenthrin	N.D.	1	0.859	1.01	85.90	101.00	16.16	
Boscalid	N.D.	1	0.817	0.888	81.70	88.80	8.33	
Carbaryl	N.D.	1	1.05	1.12	105.00	112.00	6.45	
Carbofuran	N.D.	1	0.941	0.987	94.10	98.70	4.77	
Chlorantraniliprole	N.D.	1	0.957	0.962	95.70	96.20	0.52	
Chlorfenapyr	N.D.	1	0.56	0.586	56.00	58.60	4.54	LR
Chlorpyrifos	N.D.	1	0.795	0.912	79.50	91.20	13.71	
Clofentezine	N.D.	1	0.676	0.786	67.60	78.60	15.05	
Cyfluthrin	N.D.	1	0.744	0.774	74.40	77.40	3.95	
Cypermethrin	N.D.	1	0.752	0.883	75.20	88.30	16.02	
Daminoside	N.D.	1	0.416	0.411	41.60	41.10	1.21	LR
Diazanone	N.D.	1	0.971	1.02	97.10	102.00	4.92	
Dichlorvos	N.D.	1	0.772	0.85	77.20	85.00	9.62	
Dimethoate	N.D.	1	0.927	0.937	92.70	93.70	1.07	
Ethoprophos	N.D.	1	0.86	0.916	86.00	91.60	6.31	
Etofenprox	N.D.	1	0.713	0.797	71.30	79.70	11.13	
Etoxazole	N.D.	1	1.14	1.2	114.00	120.00	5.13	
Fenoxycarb	N.D.	1	1.06	1.14	106.00	114.00	7.27	
Fenpyroximate	N.D.	1	0.886	0.954	88.60	95.40	7.39	
Fipronil	N.D.	1	1.15	1.15	115.00	115.00	0.00	
Flonicamid	N.D.	1	0.796	0.824	79.60	82.40	3.46	
Fludioxonil	N.D.	1	1.08	1.05	108.00	105.00	2.82	
Hexythiazox	N.D.	1	0.598	0.695	59.80	69.50	15.00	LR
Imazalil	N.D.	1	1.11	1.1	111.00	110.00	0.90	
Imidacloprid	N.D.	1	0.851	0.917	85.10	91.70	7.47	
Kresoxim-methyl	N.D.	1	0.995	0.998	99.50	99.80	0.30	
Malathion	N.D.	1	1.13	1.08	113.00	108.00	4.52	
Metalaxyl	N.D.	1	1.11	1.15	111.00	115.00	3.54	
Methiocarb	N.D.	1	0.885	0.921	88.50	92.10	3.99	
Methomyl	N.D.	1	1	1.01	100.00	101.00	1.00	
MGK-264	N.D.	1	0.545	0.558	54.50	55.80	2.36	LR
Myclobutanil	N.D.	1	0.916	0.982	91.60	98.20	6.95	
Naled (dibrom)	N.D.	1	0.157	0.162	15.70	16.20	3.13	LR
Oxamyl	N.D.	1	0.902	0.932	90.20	93.20	3.27	
Paclobutrazol	N.D.	1	1.04	1.06	104.00	106.00	1.90	

Parathion-methyl	N.D.	1	0.965	0.898	96.50	89.80	7.19	
Permethrins	N.D.	1	0.635	0.759	63.50	75.90	17.79	
Phosmet	N.D.	1	0.933	0.938	93.30	93.80	0.53	
Piperonyl butoxide	N.D.	1	1.06	1.09	106.00	109.00	2.79	
Prallethrin	N.D.	1	0.9	0.925	90.00	92.50	2.74	
Propiconazole	N.D.	1	0.847	0.89	84.70	89.00	4.95	
Propoxur	N.D.	1	0.943	0.976	94.30	97.60	3.44	
Pyrethrin	N.D.	0.65	0.537	0.576	82.62	88.62	7.01	
Pyridaben	N.D.	1	0.875	0.943	87.50	94.30	7.48	
SpinosynA	N.D.	0.84	0.949	0.965	112.98	114.88	1.67	
SpinosynD	N.D.	0.16	0.192	0.192	120.00	120.00	0.00	
Spiromesifen	N.D.	1	0.0686	0.0742	6.86	7.42	7.84	LR
Spirotetramat	N.D.	1	1	0.992	100.00	99.20	0.80	
Spiroxamine	N.D.	1	1.15	1.18	115.00	118.00	2.58	
Tebuconazole	N.D.	1	0.978	0.99	97.80	99.00	1.22	
Thiacloprid	N.D.	1	0.936	0.968	93.60	96.80	3.36	
Thiamethoxam	N.D.	1	0.987	1.01	98.70	101.00	2.30	
Trifloxystrobin	N.D.	1	0.756	0.75	75.60	75.00	0.80	

N.D. = Not Detected

I = indicates that an amount of an interfering compound greater than the methods limit of detection was detected in the method blank sample. May indicate contamination of analytical system or consumables.

Q = indicates that the relative percent difference of two identically prepared Matrix Spike samples for a target analyte was greater than 30%

R = indicates compound recovery of matrix spike was outside the methods acceptable limits. (60-120%) Low recovery could indicate there is actually more compound present than detected; while high recoveries should be scrutinized for possible fails as more compound may be detected than is actually residual on the sample.

Sample was sampled and tested in accordance with the Safety Compliance Facility Sampling and Testing Information.

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Test Certificate #: 128457-001

Client Name, Sample Details
Vision Aura Hemp Oil

Sample: Full Spectrum CBD
Type: Concentrate
Method: SOP FE-44-MI

Test Conditions
Prepsheet ID#: MIHS201223
Scale: XS205-MI2
Temp: 21.2 °C
Baro PE: 978.6 hPa
Analyst: KEB
Technician: ANJ

Sample ID#: 128457
Harvest/Process Date: 12/23/2020
Serving Size (g): 1
Date Received: 12/23/2020
Test Date: 12/23/2020
Valid Through: 12/23/2021
Report Issued: 12/24/2020



Compound	MRL (µg/g)	LOD (µg/g)	Status (µg/g)	Compound	MRL (µg/g)	LOD (µg/g)	Status (µg/g)
1,2-Dichloroethane	2	1	Pass/<LOD	1,2-Dimethoxyethane	5,000	250	Not Tested
1,4-dioxane	5,000	25	Not Tested	1-Butanol	5,000	250	Not Tested
1-Pentanol	5,000	250	Not Tested	1-Propanol	5,000	250	Not Tested
2,2-Dimethylpropane (Neopentane)	750	25	Pass/<LOD	2,2-Dimethylbutane (Hexanes)	50	25	Pass/<LOD
2,3-Dimethylbutane (Hexanes)	50	25	Pass/<LOD	2-Butanol	5,000	25	Not Tested
2-Butanone (MEK)	5,000	250	Not Tested	2-Ethoxyethanol	5,000	25	Not Tested
2-Methylbutane (Isopentane)	750	25	Pass/<LOD	2-Methylpentane (Hexanes)	50	25	Pass/<LOD
2-Methylpropane (Isobutane)	800	25	Pass/<LOD	2-propanol (Isopropyl Alcohol)	500	25	Pass/<LOD
2-Propanone (Acetone)	750	25	Pass/191	3-Methylpentane (Hexanes)	50	25	Pass/<LOD
Acetonitrile	60	30	Pass/<LOD	Benzene	1	1	Pass/<LOD
Butane	800	25	Pass/<LOD	Chloroform	2	1	Pass/<LOD
Cumene	5,000	25	Not Tested	Cyclohexane	50	25	Not Tested
Dichloromethane	125	25	Pass/<LOD	Dimethylsulfoxide (DMSO)	5,000	250	Not Tested
Ethanol	1,000	50	Pass/<LOD	Ethyl acetate	400	25	Pass/<LOD
Ethyl ether	500	25	Pass/<LOD	Ethylene glycol	5,000	25	Not Tested
Ethylene oxide	5	3	Pass/<LOD	Heptane	500	25	Pass/<LOD
Hexane	50	25	Pass/<LOD	Isopropyl acetate	5,000	25	Not Tested
Methanol	250	125	Pass/<LOD	Naptha	400	100	Not Tested
N,N-Dimethylacetamide	5,000	250	Not Tested	N,N-Dimethylformamide (DMF)	5,000	250	Not Tested
Pentane	750	25	Pass/<LOD	Petroleum Ether	400	100	Not Tested
Propane	2,100	25	Pass/<LOD	Pyridine	5,000	250	Not Tested
Sulfolane	5,000	25	Not Tested	Tetrahydrofuran (THF)	5,000	25	Not Tested
Toluene	150	25	Pass/<LOD	Trichloroethylene	25	25	Pass/<LOD
Xylenes*	150	25	Pass/<LOD				

Total Butanes = 0 ug/g (PASS); Total Pentanes = 0 ug/g (PASS); Total Hexanes = 0 ug/g (PASS)

* Xylenes are reported as the sum of o-xylene, m-xylene, p-xylene, and ethylbenzene

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Method: SOP FE-44-MI

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Temp: 21.2 °C
Baro PE: 978.6 hPa
Analyst: KEB
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Target Compound Name	Method Blank (µg/g)	QC Spike (µg/g)	Matrix Spike (µg/g)	Matrix Spike Duplicate (µg/g)	MS Recovery (%)	MSD Recovery (%)	Relative Percent Difference (%)	QC Flag
1,2-Dichloroethane	N.D.	250	223	223	89.20	89.20	0.00	
n-Propane	N.D.	250	108	109	43.20	43.60	0.92	LR
Isobutane	N.D.	250	148	148	59.20	59.20	0.00	LR
Methanol	N.D.	250	204	208	81.60	83.20	1.94	
n-Butane	N.D.	250	165	165	66.00	66.00	0.00	LR
2,2-Dimethylpropane	N.D.	250	179	180	71.60	72.00	0.56	
Ethanol	N.D.	250	199	208	79.60	83.20	4.42	
Isopentane	N.D.	250	189	185	75.60	74.00	2.14	
Acetonitrile	N.D.	250	196	202	78.40	80.80	3.02	
Diethyl ether	N.D.	250	209	206	83.60	82.40	1.45	
2-propanone	30.28	250	195	202	78.00	80.80	3.53	I
2-propanol	N.D.	250	198	204	79.20	81.60	2.99	
n-Pentane	N.D.	250	191	187	76.40	74.80	2.12	
2,2-Dimethylbutane	N.D.	250	233	228	93.20	91.20	2.17	
2,3-Dimethylbutane	N.D.	250	199	197	79.60	78.80	1.01	
Methylene chloride	N.D.	250	200	205	80.00	82.00	2.47	
2-Methylpentane	N.D.	250	203	194	81.20	77.60	4.53	
3-Methylpentane	N.D.	250	205	200	82.00	80.00	2.47	
2-Butanol	N.D.	0	0	0	0.00	0.00	0.00	
n-Hexane	N.D.	250	203	202	81.20	80.80	0.49	
Ethyl acetate	N.D.	250	202	204	80.80	81.60	0.99	
Tetrahydrofuran	N.D.	0	0	0	0.00	0.00	0.00	
Trichloroethylene	N.D.	250	212	209	84.80	83.60	1.43	
Isopropyl acetate	N.D.	0	0	0	0.00	0.00	0.00	
Benzene	N.D.	250	213	211	85.20	84.40	0.94	
Chloroform	N.D.	250	216	219	86.40	87.60	1.38	
Cyclohexane	N.D.	0	0	0	0.00	0.00	0.00	
Ethylene glycol	N.D.	0	0	0	0.00	0.00	0.00	
Heptane	N.D.	250	208	204	83.20	81.60	1.94	
2-Ethoxyethanol	N.D.	0	0	0	0.00	0.00	0.00	
1,4-Dioxane	N.D.	0	0	0	0.00	0.00	0.00	
Pyridine	N.D.	0	0	0	0.00	0.00	0.00	
Toluene	N.D.	250	218	219	87.20	87.60	0.46	
Chlorobenzene	N.D.	0	0	0	0.00	0.00	0.00	
Ethylbenzene	N.D.	0	0	0	0.00	0.00	0.00	
m-Xylene/p-Xylene	N.D.	500	464	459	92.80	91.80	1.08	
o-Xylene	N.D.	250	234	227	93.60	90.80	3.04	
Cumene	N.D.	0	0	0	0.00	0.00	0.00	
Ethylene oxide	N.D.	250	274	282	109.60	112.80	2.88	

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Scale: XS205-MI2

Temp: 21.2 °C

Baro Pressure: 978.6 hPa

Analyst: KEB

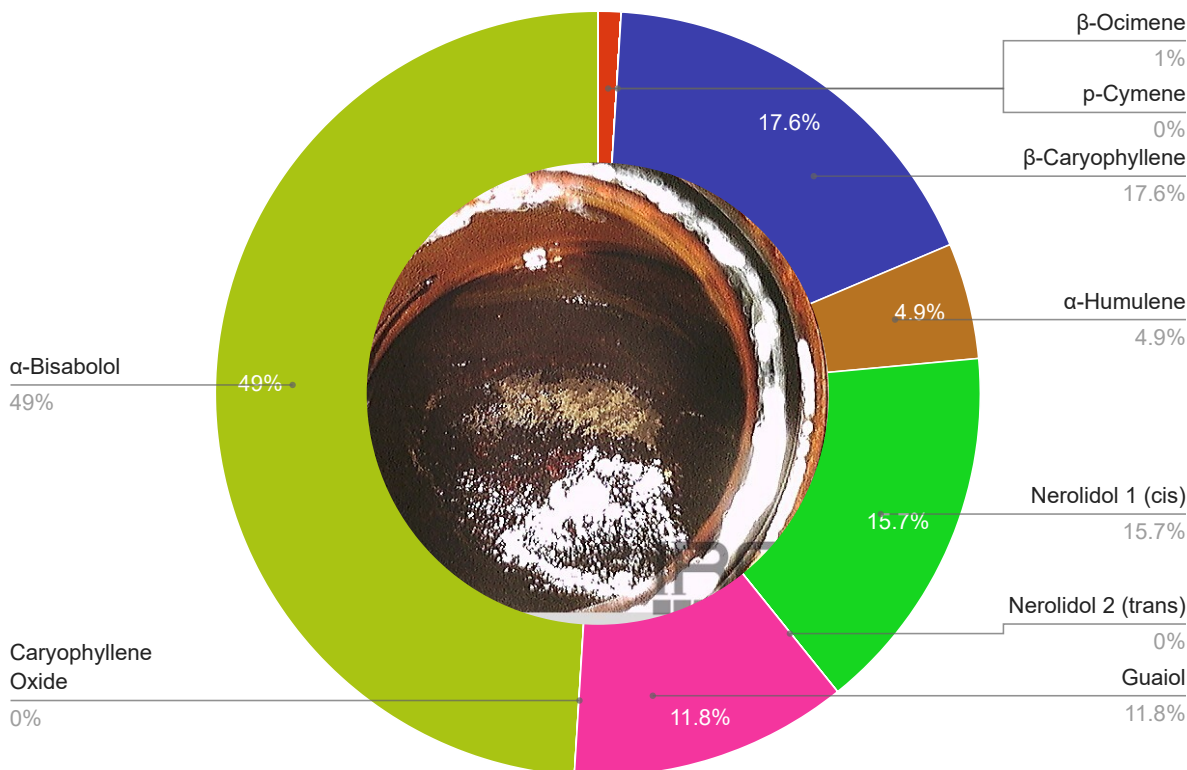
Technician: ANJ

Sample ID#: 128457

Harvest/Process Date: 12/23/2020

Serving Size (g): 1

Date Received: 12/23/2020



α-Pinene (0.00%)

γ-Terpinene (0.00%)

Δ3-Carene (0.00%)

Isopulegol (0.00%)

α-Humulene (0.05%)

α-Bisabolol (0.50%)

β-Ocimene (0.01%)

Myrcene (0.00%)

Fenchone (0.00%)

Limonene (0.00%)

Nerolidol 1 (cis) (0.16%)

α-Phellandrene (0.00%)

Camphene (0.00%)

α-Terpinolene (0.00%)

α-Terpinene (0.00%)

Geraniol (0.00%)

Nerolidol 2 (trans) (0.00%)

α-Terpineol (0.00%)

Eucalyptol (1,8-Cineol) (0.00%)

β-Pinene (0.00%)

endo-Fenchol (0.00%)

p-Cymene (0.00%)

Guaiol (0.12%)

Valencene (0.00%)

Sabinene (0.00%)

Linalool (0.00%)

α-Ocimene (0.00%)

β-Caryophyllene (0.18%)

Caryophyllene Oxide (0.00%)

Predominant Terpenes

0.50% α-Bisabolol

0.16% Nerolidol 1 (cis)

0.05% α-Humulene

Fruity, nutty, coconut

Floral, green, citrus, woody

Woody, oceanic-watery, spicy clove

0.18% β-Caryophyllene

0.12% Guaiol

0.01% β-Ocimene

Sweet, woody, spicy, clove

Mild, guaiacwood, tea, rose

Citrus, tropical, woody, green

Total: 1.020%

Results should be interpreted as qualitative only (present/absent). Quantitative values are only estimations for the purpose of research and development, and are not included in Iron Laboratories accredited scope. Eucalyptol, Sabinene, Fenchone, endo-Fenchol, Caryophyllene Oxide, α-Phellandrene, α-Terpineol, and Valencene are not currently tested for by Iron Laboratories Michigan.

Value in parenthesis indicates percentage of terpene present in the total sample (weight percentage, wt/wt%).

Value in doughnut slice indicates individual terpene abundance with respect to the total terpenes detected.

These results should be used for research and development or quality control purposes only. If applicable, results for caregivers apply to the sample as received.

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Katrina Barnes
Katrina Barnes, Lab Manager



Mac Hyman
Mackenzie E. Hyman, Quality Manager

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