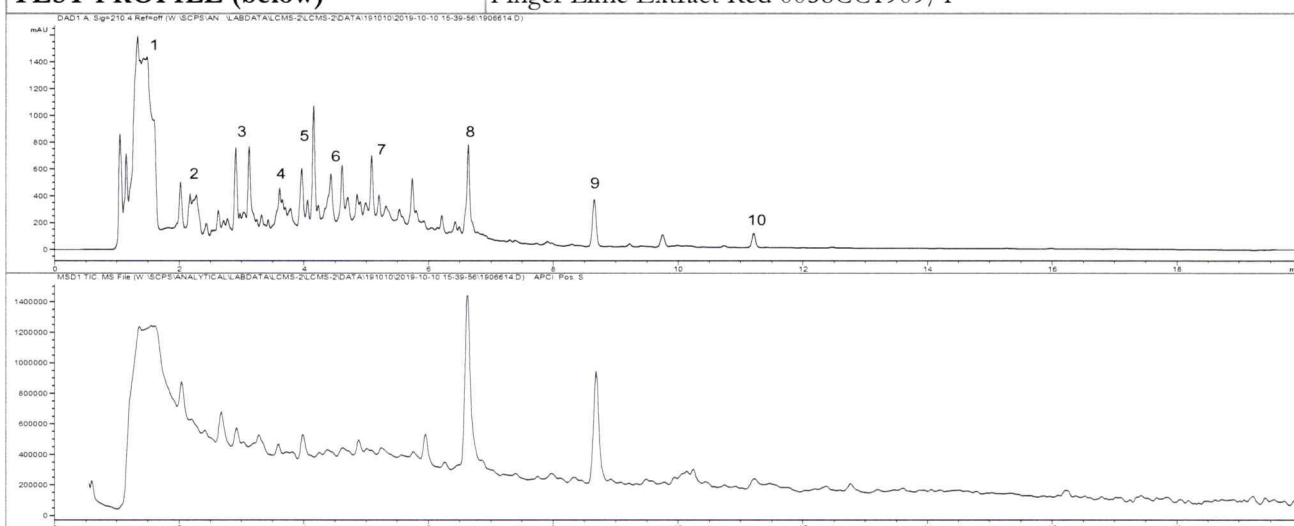


### CERTIFICATE OF ANALYSIS

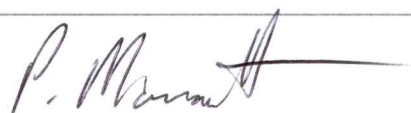
<b>SAMPLE NAME</b>		Finger Lime Extract Red	
<b>FORM</b>		Liquid	
<b>CUSTOMER NAME</b>		Plant Extracts	
<b>CERTIFICATION DATE</b>		31 October 2019	
<b>CUSTOMER REFERENCE</b>		0058CC1909/1	
<b>ARL JOB #</b>	A192425	<b>LAB REF. #</b>	ARL1906614
<b>ANALYSIS</b>	LCMS Compositional analysis	<b>METHOD</b>	ARL-TM125
<b>TEST PROFILE (below)</b>		Finger Lime Extract Red 0058CC1909/1	


**TABLE 1. PEAK IDENTIFICATION**

Peak #	RT (min)	Fragment ions [M+H]	Tentative ID (MW)
1	1.0 - 1.5	130, 216	amino acids, amines
2	2.0, 2.3	188, 205	amino acid (tryptophan), phenolic
3	2.9, 3.2	193, 231	phenolic acid, flavonoid derivative
4	3.6, 3.7	317, 449	flavonoid derivatives
5	3.9, 4.2	579, 433	flavone C-glycosides
6	4.5, 4.7	303, 317, 347, 463	flavone glycoside, phenolic derivative
7	5.1	347, 509	flavonone glycoside
8	6.7	231, 317, 335, 352, 577, 603	mixed peak - benzopyranone derivative (byangelicin), flavonone derivative
9	8.7	217, 247	benzopyranone (bergapten) derivative
10	11.2	413, 471, 531, 545	limonoid derivative

**COMMENTS**

The HPLC-MS profile of the test sample is given above with some major components from the plant extracts indicated. The major peaks identified are a range of amine or amino acids, phenolic acid, flavonoid O and C-glycosides, methoxy-benzopyranone derivatives including bergapten derivatives. Spectral data in support of peak identification is attached.

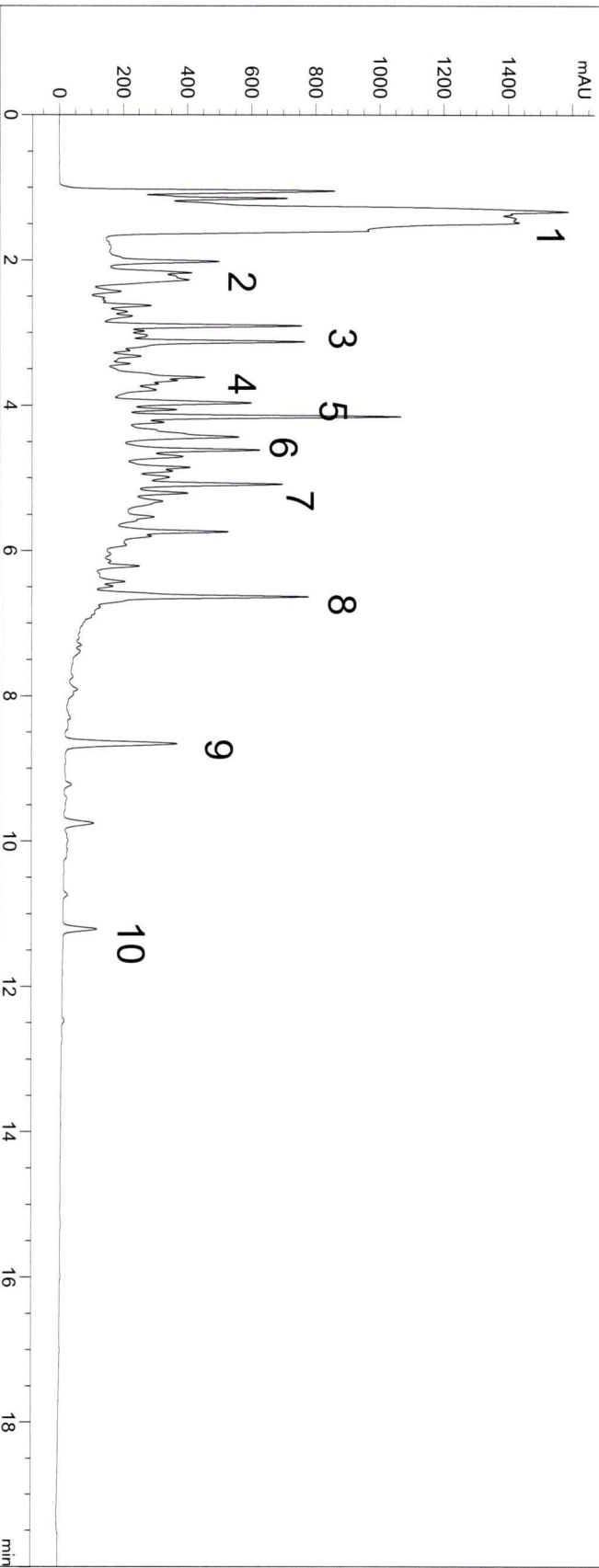


Peter Mouatt  
**SENIOR ANALYTICAL OFFICER**

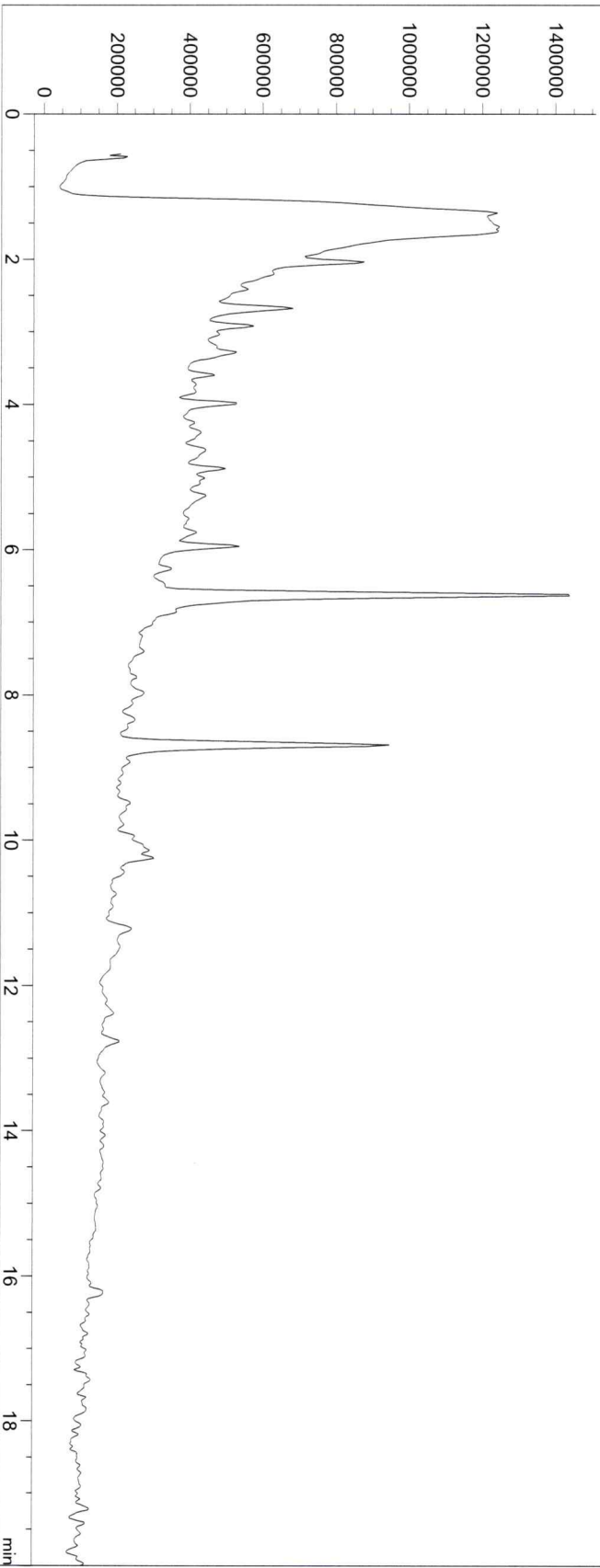

Ashley Dowell  
**MANAGER - ARL**

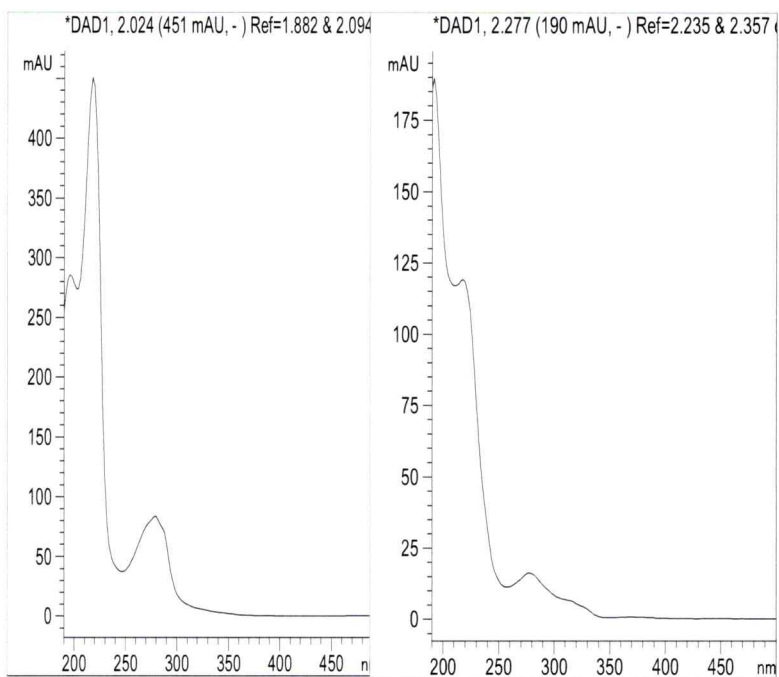
Current Chromatogram(s)

DAD1 A, Sig=210,4 Ref=off (W:\ISCPSPAN...LABDATA\LCMS-2\LCMS-2\DATA\191010\2019-10-10 15-39-56\1906614.D)

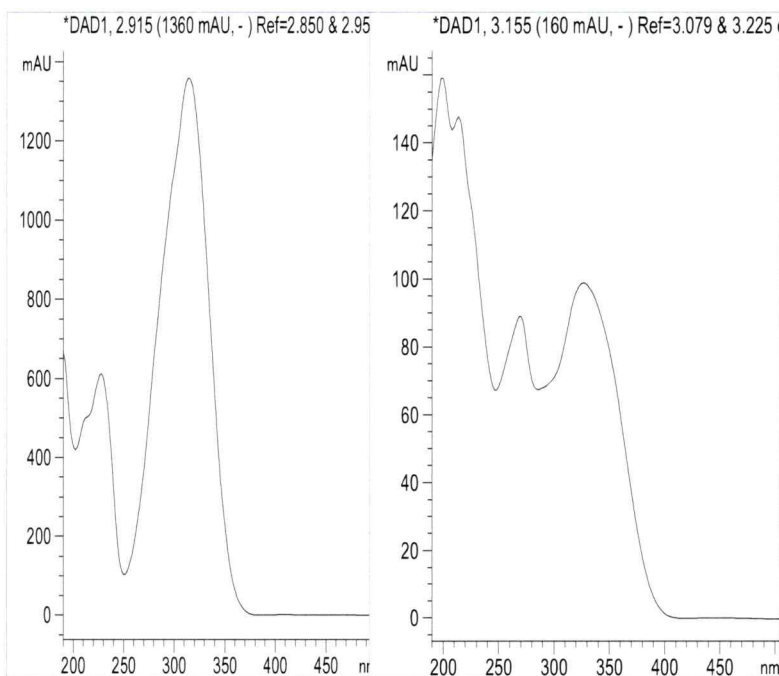


MSD1 TIC, MS File (W:\ISCPSPANALYTICAL\LABDATA\LCMS-2\LCMS-2\DATA\191010\2019-10-10 15-39-56\1906614.D) APCI, Pos, S

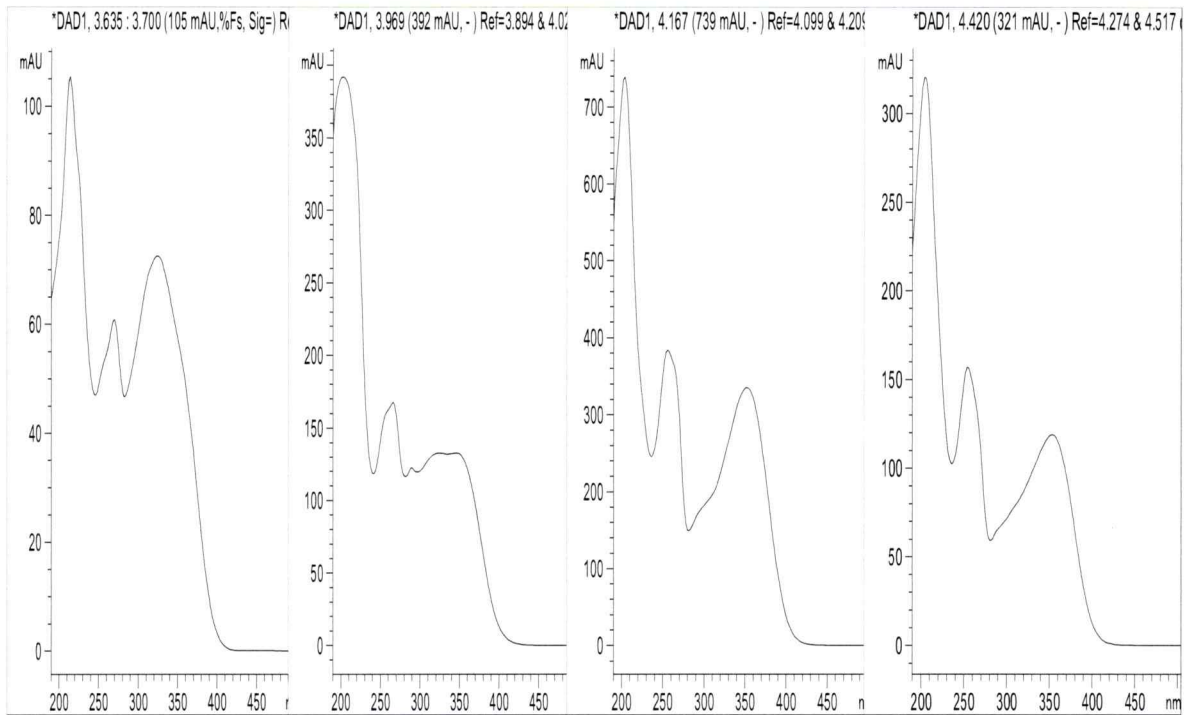




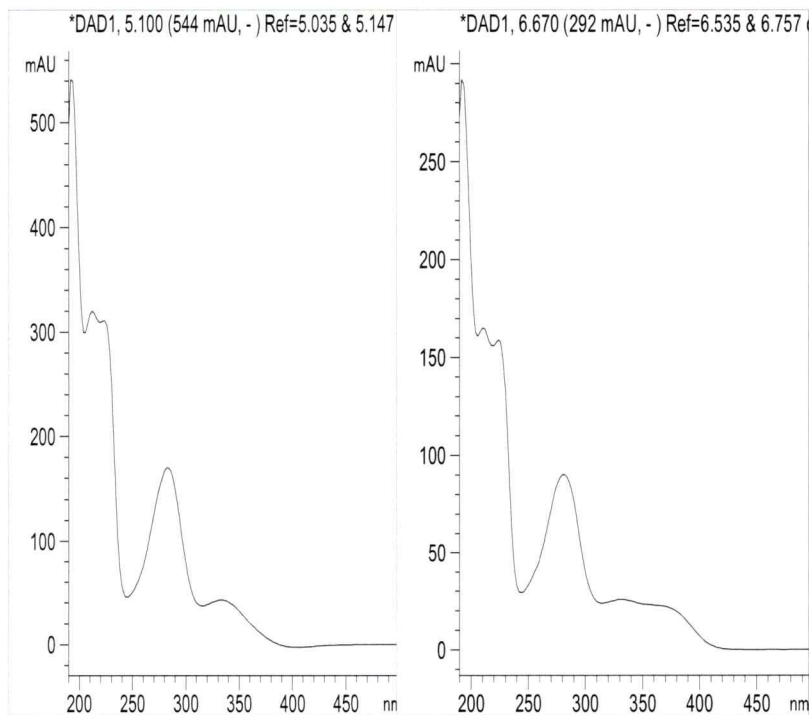
**Figure 1.** UV-Vis spectra of peaks #2 identified as amino acid tryptophan and a phenolic respectively based on UV-Vis and MS spectra



**Figure 2.** UV-Vis spectra of peaks #3 identified as phenolic acid and flavone derivative respectively based on characteristic UV-Vis spectra



**Figure 3.** UV-Vis spectra of peaks #4, #5 and #6 identified as flavone glycosides based on characteristic UV-Vis spectra



**Figure 4.** UV-Vis spectra of peaks #7 and #8 identified as flavonone derivatives based on characteristic UV-Vis spectra and absorption maxima  $\sim 280\text{nm}$