

Technical documentation

Product name:	qRE Ribes nigrum L., leaves
Substance:	Ribes nigrum L., leaves dry extract
Plant source common names:	en: Blackcurrant; fr: Cassis
Reference:	E0075
Packaging:	100 mg in a 1.5 ml borosilicate amber vial
Storage conditions:	Keep container closed. Protect from light and moisture. Keep inferior to -15 °C.
Retest:	12 months

Botanical identification of plant source

Plants in our botanical garden are identified and a herbal voucher is prepared by an expert botanist. Each batch collected for extraction is verified and identified.

Reference: Flora Europaea, Cambridge University Press, 1993, Vol 1, p 383

Method of production of dry extract

Whole plant or plant parts are collected, freeze-dried and coarsely ground. Extraction is performed by maceration in 50 % (v/v) aqueous ethanol for 48 hours at room temperature. Ethanol is then evaporated under reduced pressure at less than 40 °C and the aqueous residue is freeze-dried.

Residual water content measurement is done by Karl Fischer titration.

Organoleptic characteristics of dry extract

Colour: Light brown

Odour: Non characteristic

Form: Fine powder

Recommended methods for use

Weight a precise weight of qRE and solubilise in the recommended solvent at the concentration indicated in the HPLC or HPTLC method described in this document.

Sonicate for 90 seconds (70 W).

Filter on a 0.45 µm PVDF membrane and put the resulting solution into HPLC dispenser or apply on the HPTLC plate.

Dose and analyse your extract with qRExtract using the HPLC / HPTLC methods described in this document or using your own methods.

HPTLC

Detection of chlorogenic acid, quercetin-3-glucoside and rutin

Layer: 10 × 10 cm HPTLC Nano-Sil-20 UV 254 (Carl Roth ref. N084.1)

Thin layer conditionnement: 1 h at room temperature and 33 % relative humidity

Elution solvent:

Elution solvent compound	Volume (ml)
ethyl acetate	80
H ₂ O	10
acetic acid	10

Initial spot volume and concentration:

quercetin-3-glucoside: 1 µl of a 0.08 % (w/v) solution in ethanol ethanol 96 %
 rutin: 3 µl of a 0.02% (w/v) solution in methanol
 qRE: 5 µl of a 1.6 % (w/v) solution in 50 % (v/v) aqueous ethanol
 chlorogenic acid: 1 µl of a 0.02 % (w/v) solution in 50 % (v/v) aqueous ethanol

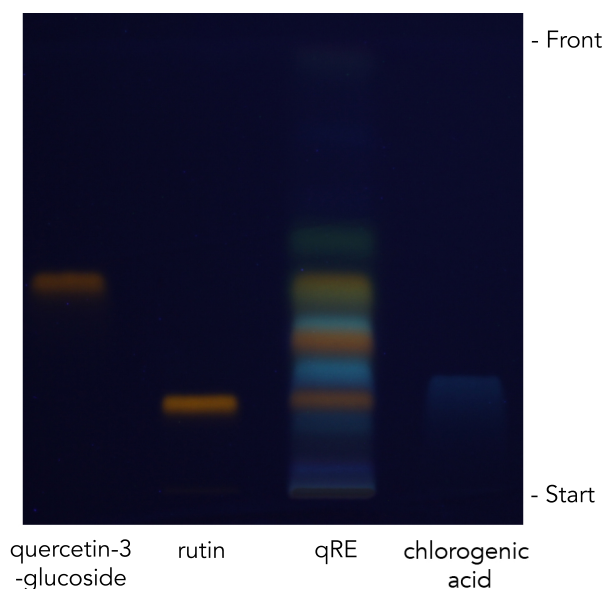
Reagent mixture:

Natural products - polyethylene glycol reagent (NP/PEG)

Preparation: Dissolve 0.25 g of diphenylboric acid 2-aminoethylester and 1.25 g of polyethylene glycol 400 in 25 mL of methanol.

Dip the plate in the reagent mixture and dry for 15 minutes at room temperature.

Expose to UV light at 365 nm.

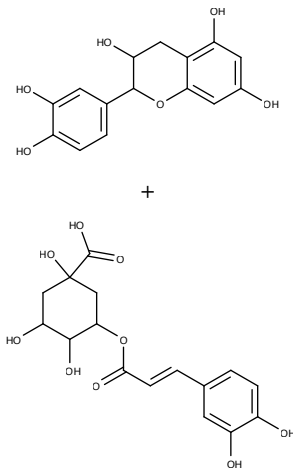


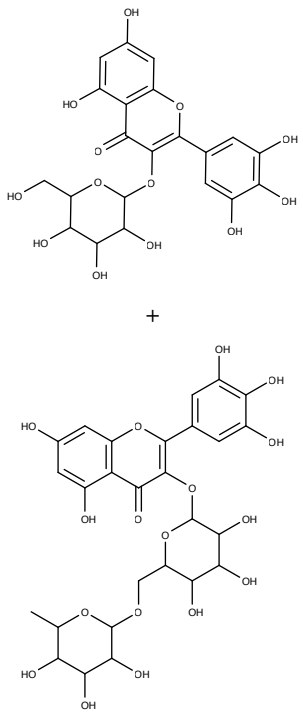
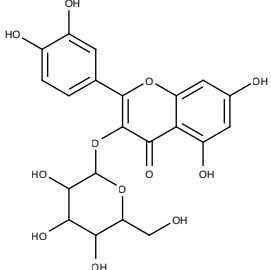
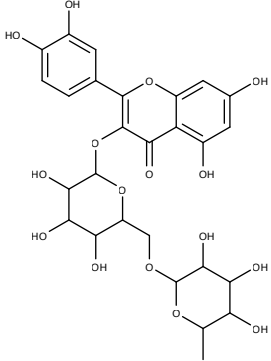
HPLC

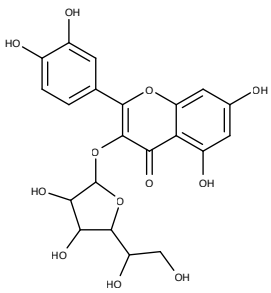
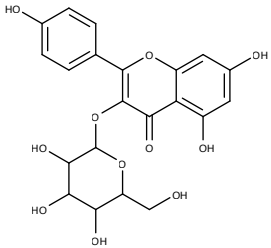
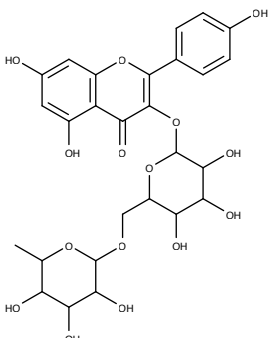
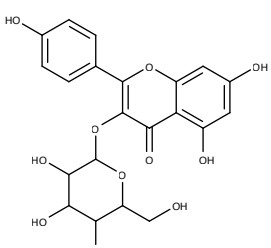
Precolumn: Ascentis® Express C18 0.5 cm × 3.0 mm 2.7 μm
Column: Ascentis® Express C18 15 cm × 3.0 mm 2.7 μm
Sample: 8 μl 3.0 % qRE® (w/v) solution in 25 % (v/v) aqueous ethanol
Flow: 0.45 ml/min
Temperature: 25 °C
Mobile phase: A: 0.1 % formic acid (v/v) in water
 B: 0.1 % formic acid (v/v) in acetonitrile
Detection: Diode Array Detector, 350 nm
Gradient:

Time (mn)	A %	B %
0	97	3
50	87	13
110	87	13
118	86.4	13.6

Quantified substances

Compound	CAS No	2D Structure	Peak No
Epicatechin + Chlorogenic acid	490-46-0 + 327-97-9	 <p>The image shows the chemical structures of Epicatechin and Chlorogenic acid. Epicatechin is a flavan-3-ol consisting of a flavan-3-ol core with a catechol B-ring and a galloyl C-ring. Chlorogenic acid is a hydroxycinnamic acid derivative consisting of a 3,4-dihydroxyphenyl ring connected via an ester linkage to a 3,4-dihydroxyphenyl ring.</p>	2

Compound	CAS No	2D Structure	Peak No
Myricetin glucoside + myricetin rutinoside	19833-12-6 + NA	 <p>The image shows two chemical structures. The top structure is Myricetin glucoside, consisting of a myricetin aglycone (a flavan-3-ol with three hydroxyl groups on the A-ring and two on the C-ring) linked to a glucose molecule at the 3-position. The bottom structure is Myricetin rutinoside, where the myricetin aglycone is linked to a rutinose molecule (a disaccharide of glucose and rhamnose) at the 3-position. A plus sign is placed between the two structures.</p>	3
Quercetin-3-O-galactoside	482-36-0	 <p>The image shows the chemical structure of Quercetin-3-O-galactoside. It features a quercetin aglycone (a flavonol with two hydroxyl groups on the A-ring and three on the C-ring) linked to a galactose molecule at the 3-position.</p>	6
Quercetin-3-O-rutinoside	153-18-4	 <p>The image shows the chemical structure of Quercetin-3-O-rutinoside. It features a quercetin aglycone linked to a rutinose molecule (a disaccharide of glucose and rhamnose) at the 3-position.</p>	7

Compound	CAS No	2D Structure	Peak No
Quercetin-3-O-glucoside	482-35-9		8
Kaempferol-3-glucoside isomer	0480-10-4		9
Kaempferol-3-O-rutinoside	17650-84-9		12
Kaempferol-3-glucoside isomer	0480-10-4		13
Unknown	NA	NA	1, 4, 5, 10, 11, 14