

Thermo EASY-LC method print for Thermo

Sample pickup:

Volume [µl] : 4.00  
Flow [µl / min] : 10.00

Sample loading:

Volume [µl] : 10.00  
Flow [µl / min] : (unspecified)  
Max. pressure [Bar] : 750.00

Gradient:

Time [mm:ss]	Duration [mm:ss]	Flow [nl/min]	Mixture [%B]
00:00	00:00	250.00	2.00
05:00	05:00	250.00	8.00
80:00	75:00	250.00	28.00
90:00	10:00	250.00	44.00
95:00	05:00	250.00	100.00
98:00	03:00	300.00	100.00
101:00	03:00	300.00	2.00
104:00	03:00	300.00	2.00
107:00	03:00	300.00	100.00
110:00	03:00	300.00	100.00

Pre-column equilibration:

Volume [µl] : 0.00  
Flow [µl / min] : (unspecified)  
Max. pressure [Bar] : 750.00

Analytical column equilibration:

Volume [µl] : 3.00  
Flow [µl / min] : (unspecified)  
Max. pressure [Bar] : 750.00

Autosampler wash:

Flush volume [µl] : 100.00

## Method Summary

### Method Settings

Application Mode: **Peptide**  
Method Duration (min): **110**

### Global Parameters

#### Ion Source

Ion Source Type: **NSI**  
Spray Voltage: **Static**  
Positive Ion (V): **2000**  
Negative Ion (V): **600**  
Sweep Gas (Arb): **0**  
Ion Transfer Tube Temp (°C): **280**  
Use Ion Source Settings from Tune: **False**  
FAIMS Mode: **Not Installed**

#### MS Global Settings

Infusion Mode: **Liquid Chromatography**  
Expected LC Peak Width (s): **20**  
Advanced Peak Determination: **True**  
Default Charge State: **2**  
Internal Mass Calibration: **EASY-IC™**  
Mode: **Run Start**

### Experiment#1 [tSIM]

Start Time (min): **0**  
End Time (min): **110**

#### Master Scan:

#### tSIM

Multiplex Ions: **True**  
Maximum number of multiplexed ions: **6**  
Define Multiplexing Groups (MSX ID): **User-defined**  
Isolation Offset: **Off**

Orbitrap Resolution: **60000**  
RF Lens (%): **40**  
AGC Target: **Custom**  
Normalized AGC Target (%): **1000**  
Maximum Injection Time Mode: **Custom**  
Maximum Injection Time (ms): **4**  
Microscans: **1**  
Data Type: **Centroid**  
Polarity: **Positive**  
Source Fragmentation: **Disabled**  
Scan Description:  
Time Mode: **Unscheduled**

### Mass List Table

Mass List Table						
Compound	Formula	Adduct	m/z	z	MSX ID	Isolation Window (m/z)
Boxcar1		(no adduct)	413.4378	1	1	126
Boxcar2		(no adduct)	521.4869	1	1	90
Boxcar3		(no adduct)	611.0276	1	1	89
Boxcar4		(no adduct)	705.5706	1	1	100
Boxcar5		(no adduct)	823.124	1	1	135.1
Boxcar6		(no adduct)	1045.7252	1	1	310.1

### Filters:

#### MIPS

Monoisotopic peak determination: **Peptide**  
Relax restrictions when too few precursors are found: **True**

#### Intensity

Filter Type: **Intensity Threshold**  
Intensity Threshold: **8.0e3**

#### Charge State

Include charge state(s): **2-6**  
Include undetermined charge states: **False**

## Dynamic Exclusion

Dynamic Exclusion Mode: **Custom**  
Exclude after n times: **1**  
Exclusion duration (s): **30**  
Mass Tolerance: **ppm**  
Low: **10**  
High: **10**  
Exclude isotopes: **True**  
Perform dependent scan on single charge state per precursor only: **True**

## Data Dependent

Data Dependent Mode: **Number of Scans**  
Number of Dependent Scans: **20**

## Scan Event Type 1:

Scan:

ddMS<sup>2</sup>

Multiplex Ions: **False**  
Isolation Window (m/z): **2**  
Isolation Offset: **Off**  
Collision Energy Type: **Normalized**  
HCD Collision Energy (%): **27**  
Orbitrap Resolution: **15000**  
TurboTMT: **Off**  
Scan Range Mode: **Define First Mass**  
First Mass (m/z): **140**  
AGC Target: **Custom**  
Normalized AGC Target (%): **50**  
Maximum Injection Time Mode: **Auto**  
Microscans: **1**  
Data Type: **Centroid**  
Scan Description: