Analysis Report

Bruker IVDr Quantification in URine B.I.Quant-UR b™

Sample ID:	ALZ_Urine_Rack01_RCM_221214_expno570.100000.10r

Measuring Date:	24-Dec-2014 03:29:18
Reporting Date:	12-Dec-2020 11:13:55, 7 page(s), Version 1.1.0
Quantification Method Version:	Quant-UR B.1.1.0

Disclaimer

RESEARCH USE ONLY: This is no clinical diagnostic analysis report. Must not be used for clinical (medical or IVD) diagnosis or for patient management! Additional concentration range information (95% range) provided numerically or graphically in this report must not be used for clinical diagnostic interpretation.

Application of B.I.Quant-UR B 1.1.0 requires use of Bruker's B.I.Methods SOP for urine.

Summary

The following metabolites were found with concentrations outside the 95% range of Bruker Quant-UR B.1.1.0 urine metabolite concentration database:

Amino acids and derivatives: Betaine (< 7 mmol/mol Crea).

Further detailed information is provided on the following pages.

Contents

1 Creatinine

Compound	Conc.	LOD	r	ρ	Δ	95% Range[*]
	mmol/L	mmol/L	mmol/L	%	mmol/L	mmol/L
Creatinine	14	0.3	14.16	100 🟢	0.328	1 - 19

[*] Gray horizontal boxes represent 95% concentration range, black vertical lines represent sample value.

2 Amines and derivatives

Compound	Conc.	Conc.	LOD	r	ρ	Δ	95% Range[*]
	mmol/L	$\frac{mmol}{mol\ Crea}$	$\frac{mmol}{mol\ Crea}$	mmol/L	%	mmol/L	$\frac{mmol}{mol\ Crea}$
Dimethylamine	< 0.43	< 31	31	0.429	100 🟢	0.016	≤ 54
Trimethylamine	< 0.03	< 2	2	0.002	10 ⚪	0.002	≤ 3

[*] Gray horizontal boxes represent 95% concentration range, black vertical lines represent sample value.

3 Amino acids and derivatives

Compound	Conc.	Conc.	LOD	r	ρ	Δ	95% Range[*]
	mmol/L	$\frac{mmol}{mol\ Crea}$	$\frac{mmol}{mol\ Crea}$	mmol/L	%	mmol/L	$\frac{mmol}{mol\ Crea}$
1-Methylhistidine	< 0.21	< 15	15	0.000	0 ⚪	0.169	≤ 15
2-Furoylglycine	< 0.55	< 39	39	0.252	79 ⚪	0.051	≤ 40
4-Aminobutyric acid	< 0.28	< 20	20	0.188	58 ⚪	0.269	≤ 20
Alanine	0.19	13	10	0.186	99 🟢	0.025	11 - 72
Arginine	< 11	< 750	750	0.881	39 ⚪	1.188	≤ 750
Betaine	< 0.10	< 7	7	0.097	100 🟢	0.042	9 - 78
Creatine	< 0.71	< 50	50	0.002	100 🟢	0.328	≤ 280
Glycine	0.69	49	34	0.689	100 🟢	0.051	38 - 440
Guanidinoacetic acid	< 1.5	< 100	100	0.405	76 ⚪	0.217	≤ 140
Methionine	< 0.25	< 18	18	0.000	0 ⚪	0.600	≤ 18
N,N-Dimethylglycine	0.07	5	5	0.074	67 ⚪	0.033	≤ 15
Sarcosine	< 0.03	< 2	2	0.015	0 ⚪	0.023	≤ 7
Taurine	< 2.0	< 140	140	0.689	98 🟢	0.235	≤ 170
Valine	0.05	4	2	0.051	88 🟡	0.022	≤ 7

[*] Gray horizontal boxes represent 95% concentration range, black vertical lines represent sample value.

4 Benzene and substituted derivatives

Compound	Conc. mmol/L	Conc. $\frac{mmol}{mol\ Crea}$	LOD $\frac{mmol}{mol\ Crea}$	r mmol/L	ρ %	Δ mmol/L	95% Range[*] $\frac{mmol}{mol\ Crea}$
Benzoic acid	< 0.14	< 10	10	0.000	0 ○	0.088	≤ 10
D-Mandelic acid	< 0.03	< 2	2	0.000	0 ○	0.010	2 - 17
Hippuric acid	< 2.4	< 170	170	1.588	99 ●	0.143	≤ 660

[*] Gray horizontal boxes represent 95% concentration range, black vertical lines represent sample value.

5 Carboxylic acids

Compound	Conc. mmol/L	Conc. $\frac{mmol}{mol\ Crea}$	LOD $\frac{mmol}{mol\ Crea}$	r mmol/L	ρ %	Δ mmol/L	95% Range[*] $\frac{mmol}{mol\ Crea}$
Acetic acid	< 0.07	< 5	5	0.049	91 ●	0.020	≤ 51
Citric acid	1.5	110	40	1.504	99 ●	0.335	≤ 700
Formic acid	< 0.13	< 10	10	0.097	100 ●	0.004	≤ 43
Fumaric acid	< 0.03	< 2	2	0.003	96 ●	0.001	≤ 3
Imidazole	< 0.68	< 48	48	0.511	96 ●	0.090	≤ 48
Lactic acid	< 0.69	< 49	49	0.142	88 ●	0.087	≤ 110
Proline betaine	1.5	110	25	1.546	100 ●	0.066	≤ 280
Succinic acid	< 0.07	< 5	5	0.057	95 ●	0.011	≤ 39
Tartaric acid	< 0.07	< 5	5	0.026	49 ○	0.030	≤ 110
Trigonelline	< 0.49	< 35	35	0.125	100 ●	0.003	≤ 67

[*] Gray horizontal boxes represent 95% concentration range, black vertical lines represent sample value.

6 Fatty acids and derivatives

Compound	Conc. mmol/L	Conc. $\frac{mmol}{mol\ Crea}$	LOD $\frac{mmol}{mol\ Crea}$	r mmol/L	ρ %	Δ mmol/L	95% Range[*] $\frac{mmol}{mol\ Crea}$
2-Methylsuccinic acid	< 0.68	< 48	48	0.000	0 ○	0.365	≤ 48

[*] Gray horizontal boxes represent 95% concentration range, black vertical lines represent sample value.

7 Keto acids and derivatives

Compound	Conc. mmol/L	Conc. $\frac{mmol}{mol\ Crea}$	LOD $\frac{mmol}{mol\ Crea}$	r mmol/L	ρ %	Δ mmol/L	95% Range[*] $\frac{mmol}{mol\ Crea}$
2-Oxoglutaric acid	< 1.3	< 92	92	0.091	71 ○	0.474	≤ 92
3-Hydroxybutyric acid	< 1.5	< 100	100	0.000	0 ○	1.962	≤ 100
Acetoacetic acid	< 0.20	< 14	14	0.124	84 ○	0.095	≤ 30
Acetone	0.03	2	2	0.027	98 ●	0.009	≤ 7
Oxaloacetic acid	< 0.24	< 17	17	0.211	82 ○	0.145	≤ 66
Pyruvic acid	< 0.13	< 9	9	0.057	86 ●	0.017	≤ 13

[*] Gray horizontal boxes represent 95% concentration range, black vertical lines represent sample value.

8 Purine, Pyridine and Pyrimidine derivatives

Compound	Conc. mmol/L	Conc. $\frac{mmol}{mol\ Crea}$	LOD $\frac{mmol}{mol\ Crea}$	r mmol/L	ρ %	Δ mmol/L	95% Range[*] $\frac{mmol}{mol\ Crea}$
1-Methyladenosine	< 0.07	< 5	5	0.000	0 ○	0.181	≤ 5
1-Methylnicotinamide	< 0.45	< 32	32	0.072	99 ●	0.006	≤ 32
Adenosine	< 5.5	< 390	390	0.000	0 ○	2.382	≤ 390
Allantoin	< 0.24	< 17	17	0.064	67 ○	0.036	≤ 47
Allopurinol	< 0.14	< 10	10	0.095	70 ○	0.084	≤ 11
Caffeine	< 0.64	< 45	45	0.209	87 ●	0.253	≤ 61
Inosine	< 0.27	< 19	19	0.135	51 ○	0.100	≤ 19

[*] Gray horizontal boxes represent 95% concentration range, black vertical lines represent sample value.

9 Sugars and derivatives

Compound	Conc. mmol/L	Conc. $\frac{mmol}{mol\ Crea}$	LOD $\frac{mmol}{mol\ Crea}$	r mmol/L	ρ %	Δ mmol/L	95% Range[*] $\frac{mmol}{mol\ Crea}$
D-Galactose	< 0.61	< 43	43	0.027	0 ○	0.024	≤ 44
D-Glucose	< 0.48	< 34	34	0.456	92 ●	0.106	≤ 140
D-Lactose	< 1.4	< 96	96	0.025	44 ○	0.049	≤ 96
D-Mannitol	< 2.6	< 180	180	0.000	0 ○	3.919	≤ 180
D-Mannose	< 0.08	< 6	6	0.000	0 ○	0.103	≤ 8
Myo-Inositol	< 63	< 4400	4400	0.000	0 ○	12.09	≤ 4400

[*] Gray horizontal boxes represent 95% concentration range, black vertical lines represent sample value.

10 Explanations

This section contains the definition of the parameters used above. In the section 10.1 a short manual, how to interpret the results, is presented. The section 10.3 contains the exact definitions of the parameters $\mathbf{r}$, ρ and $\mathbf{\Delta}$.
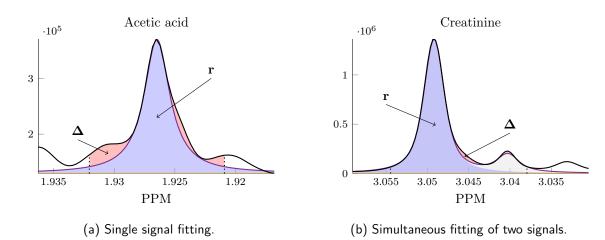
10.1 How to read the result

(a) Single signal fitting. (b) Simultaneous fitting of two signals.

Figure 1: Examples of fitting.

In the figure 1(a), the black line, the blue line and the yellow line represent the original spectrum, the calculated signal fit and its baseline, respectively.

The blue area relates to the metabolite concentration to be determined and the red area represents a residue.

In case of the signal overlap a different approach is used: two or more overlapping signals are being fitted simultaneously. The most iconic example of such signals are the ones generated by CH_3 groups of Creatinine and Creatine. In such a case, the blue line and the grey area relate the sum of all fitted signals. The blue area corresponds to the concentration of the metabolite of interest (cf. figure 1(b)).

10.2 Result parameters

a) **Conc.** is the final result concentration of the metabolite,

b) **LOD** is the *limit of detection* of the given metabolite,

c) $\mathbf{r}$ is the *raw concentration* i.e. the concentration equivalent of the resulting signal fit prior to comparing to **LOD** (relates to the blue area, cf. α)),

d) ρ is the correlation of lineshape metabolite signal with calculated fit characterizing the match between metabolite signal and fit (cf. β)). Depending on the value of ρ, the following *flag* is displayed:

- 🟢, if the correlation is 95%,

- 🟡, if the correlation is in between 85% and 95%,

- ⚪, if the correlation is less than 85%,

e) $\boldsymbol{\Delta}$ is the concentration equivalent of the difference between metabolite signal and calculated fit (residue corresponding to the the red area , cf. γ)).

10.3   Detailed definitions

Let s, f and b denote the functions describing the *raw spectra*, *fitted curve* and *(fitted) baseline* respectively. These functions are chosen such that $s \approx f + b$. Moreover, let I be a relevant PPM interval and P_N be the proton number for given metabolite/signal.

α) $\mathbf{r}$ (*raw concentration*) is defined as

$$\mathbf{r} = \frac{1}{P_N} \int_{\mathbb{R}} f(\xi)\, \mathrm{d}\xi.$$

β) ρ is the *correlation* of the functions s and $f + b$, i.e.

$$\rho = \max\left(0, \mathrm{corr}\left(\overline{s}, \overline{f+b}\right)\right),$$

where $\overline{s}$, $\overline{f+b}$ are numerical representations of the functions s and $f + b$ on sufficiently fine mesh of the interval I.

γ) $\boldsymbol{\Delta}$ is the the area between the raw signal s and the fitted data $f + b$ on the interval I expressed in the terms of the concentration, i.e.

$$\boldsymbol{\Delta} = \frac{1}{P_N} \int_{I} |s(\xi) - f(\xi) - b(\xi)|\, \mathrm{d}\xi.$$