Analysis Report

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

Sample ID: ALZ_Urine_Rack01_RCM_221214_expno620.100000.10r

Measuring Date: 24-Dec-2014 04:43:12
Reporting Date: 12-Dec-2020 15:25:58, 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),*
Carboxylic acids: *Acetic acid (150 mmol/mol Crea).*

Further detailed information is provided on the following pages.

Contents

1 Creatinine

Compound	Conc. mmol/L	LOD mmol/L	r mmol/L	ρ %	Δ mmol/L	95% Range(*) mmol/L
Creatinine	4.9	0.3	4.900	100 ●	0.134	1 - 19

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

2 Amines and derivatives

Compound	Conc. mmol/L	Conc. $\frac{\text{mmol}}{\text{mol Crea}}$	LOD $\frac{\text{mmol}}{\text{mol Crea}}$	r mmol/L	ρ %	Δ mmol/L	95% Range(*) $\frac{\text{mmol}}{\text{mol Crea}}$
Dimethylamine	0.18	38	31	0.184	100 ●	0.007	≤ 54
Trimethylamine	0.01	2	2	0.012	100 ●	0.001	≤ 3

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

3 Amino acids and derivatives

Compound	Conc. mmol/L	Conc. $\frac{\text{mmol}}{\text{mol Crea}}$	LOD $\frac{\text{mmol}}{\text{mol Crea}}$	r mmol/L	ρ %	Δ mmol/L	95% Range(*) $\frac{\text{mmol}}{\text{mol Crea}}$
1-Methylhistidine	< 0.07	< 15	15	0.027	52 ○	0.018	≤ 15
2-Furoylglycine	< 0.19	< 39	39	0.000	0 ○	0.026	≤ 40
4-Aminobutyric acid	< 0.10	< 20	20	0.000	0 ○	0.186	≤ 20
Alanine	0.08	16	10	0.076	100 ●	0.005	11 - 72
Arginine	< 3.7	< 750	750	0.298	0 ○	1.459	≤ 750
Betaine	< 0.03	< 7	7	0.008	100 ●	0.167	9 - 78
Creatine	1.00	210	50	1.035	100 ●	0.134	≤ 280
Glycine	0.27	55	34	0.268	100 ●	0.011	38 - 440
Guanidinoacetic acid	< 0.51	< 100	100	0.196	98 ●	0.021	≤ 140
Methionine	< 0.09	< 18	18	0.000	0 ○	0.145	≤ 18
N,N-Dimethylglycine	0.03	5	5	0.026	79 ○	0.008	≤ 15
Sarcosine	< 0.01	< 2	2	0.006	0 ○	0.012	≤ 7
Taurine	< 0.70	< 140	140	0.428	93 ●	0.167	≤ 170
Valine	< 0.01	< 2	2	0.005	56 ○	0.004	≤ 7

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

4 Benzene and substituted derivatives

Compound	Conc. mmol/L	Conc. $\frac{\text{mmol}}{\text{mol Crea}}$	LOD $\frac{\text{mmol}}{\text{mol Crea}}$	r mmol/L	ρ %	Δ mmol/L	95% Range(*) $\frac{\text{mmol}}{\text{mol Crea}}$
Benzoic acid	< 0.05	< 10	10	0.000	0 ○	0.029	≤ 10
D-Mandelic acid	< 0.01	< 2	2	0.000	0 ○	0.089	2 - 17
Hippuric acid	< 0.83	< 170	170	0.787	99 ●	0.063	≤ 660

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

5 Carboxylic acids

Compound	Conc. mmol/L	Conc. $\frac{\text{mmol}}{\text{mol Crea}}$	LOD $\frac{\text{mmol}}{\text{mol Crea}}$	r mmol/L	ρ %	Δ mmol/L	95% Range(*) $\frac{\text{mmol}}{\text{mol Crea}}$
Acetic acid	0.76	150	5	0.756	100 ●	0.036	≤ 51
Citric acid	1.3	270	40	1.344	100 ●	0.183	≤ 700
Formic acid	< 0.05	< 10	10	0.007	89 ●	0.003	≤ 43
Fumaric acid	< 0.01	< 2	2	0.004	92 ●	0.001	≤ 3
Imidazole	< 0.23	< 48	48	0.063	40 ○	0.083	≤ 48
Lactic acid	< 0.24	< 49	49	0.059	89 ●	0.024	≤ 110
Proline betaine	0.32	65	25	0.321	92 ●	0.055	≤ 280
Succinic acid	0.05	11	5	0.054	97 ●	0.009	≤ 39
Tartaric acid	< 0.02	< 5	5	0.012	60 ○	0.007	≤ 110
Trigonelline	< 0.17	< 35	35	0.088	100 ●	0.002	≤ 67

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

6 Fatty acids and derivatives

Compound	Conc. mmol/L	Conc. $\frac{\text{mmol}}{\text{mol Crea}}$	LOD $\frac{\text{mmol}}{\text{mol Crea}}$	r mmol/L	ρ %	Δ mmol/L	95% Range(*) $\frac{\text{mmol}}{\text{mol Crea}}$
2-Methylsuccinic acid	< 0.23	< 48	48	0.000	0 ○	0.161	≤ 48

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

7 Keto acids and derivatives

Compound	Conc. mmol/L	Conc. $\frac{\text{mmol}}{\text{mol Crea}}$	LOD $\frac{\text{mmol}}{\text{mol Crea}}$	r mmol/L	ρ %	Δ mmol/L	95% Range(*) $\frac{\text{mmol}}{\text{mol Crea}}$
2-Oxoglutaric acid	< 0.45	< 92	92	0.048	81 ○	0.112	≤ 92
3-Hydroxybutyric acid	< 0.51	< 100	100	0.000	0 ○	0.404	≤ 100
Acetoacetic acid	< 0.07	< 14	14	0.057	83 ○	0.030	≤ 30
Acetone	0.01	2	2	0.012	96 ●	0.003	≤ 7
Oxaloacetic acid	0.08	17	17	0.083	90 ●	0.048	≤ 66
Pyruvic acid	< 0.04	< 9	9	0.019	93 ●	0.004	≤ 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. $\frac{\text{mmol}}{\text{mol Crea}}$	LOD $\frac{\text{mmol}}{\text{mol Crea}}$	r mmol/L	ρ %	Δ mmol/L	95% Range(*) $\frac{\text{mmol}}{\text{mol Crea}}$
1-Methyladenosine	< 0.02	< 5	5	0.000	0 ○	0.089	≤ 5
1-Methylnicotinamide	< 0.15	< 32	32	0.056	98 ●	0.007	≤ 32
Adenosine	< 1.9	< 390	390	0.000	0 ○	0.829	≤ 390
Allantoin	< 0.08	< 17	17	0.056	99 ●	0.004	≤ 47
Allopurinol	< 0.05	< 10	10	0.016	95 ●	0.032	≤ 11
Caffeine	< 0.22	< 45	45	0.092	97 ●	0.053	≤ 61
Inosine	< 0.09	< 19	19	0.031	88 ●	0.033	≤ 19

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

9 Sugars and derivatives

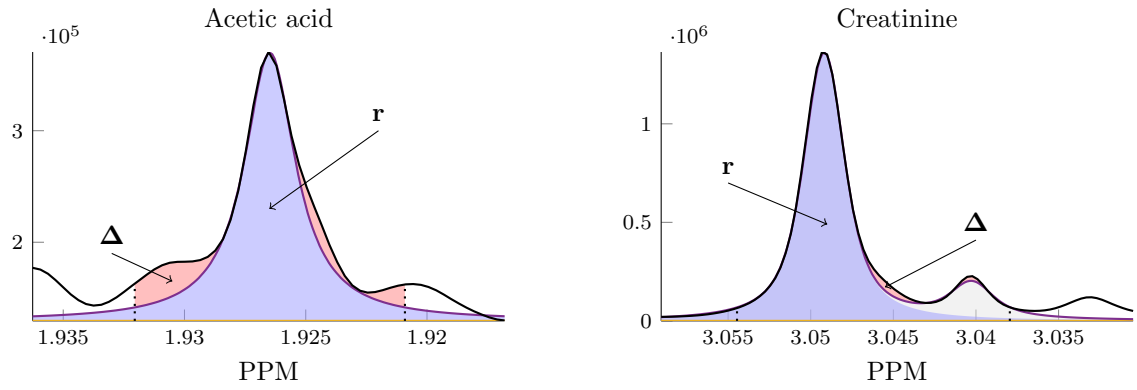
Compound	Conc. mmol/L	Conc. $\frac{\text{mmol}}{\text{mol Crea}}$	LOD $\frac{\text{mmol}}{\text{mol Crea}}$	r mmol/L	ρ %	Δ mmol/L	95% Range(*) $\frac{\text{mmol}}{\text{mol Crea}}$
D-Galactose	< 0.21	< 43	43	0.000	0 ○	0.021	≤ 44
D-Glucose	< 0.17	< 34	34	0.089	97 ●	0.011	≤ 140
D-Lactose	< 0.47	< 96	96	0.010	55 ○	0.021	≤ 96
D-Mannitol	< 0.90	< 180	180	0.000	0 ○	1.051	≤ 180
D-Mannose	< 0.03	< 6	6	0.000	0 ○	0.033	≤ 8
Myo-Inositol	< 22	< 4400	4400	0.000	0 ○	45.89	≤ 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 $\mathrm{CH_3}$ 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) **Δ** 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.

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

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) **Δ** 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.

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