

REF : AUML-XXXX

Consult Instruction for Use

APPLICATION - Mindray - BS 360 - PROPOSAL This application is intended to serve as a guide for using the referenced ELITechGroup Reagent on this instrument system only. It is recommended that the user validate this application prior to routine use.				
system only. It is recommended that the user validate this application prior to routine use.				
Define / Edite Chemestries				
Chem AUML N°			Sample Type	Serum
Chemistry URIC ACID			Print name	
Reaction Type Endpoint		Reaction Dire	ection Positive	
Pri Wave 546 nm		Sec	Wave	
Unit mg/dL			cimal 0.01]
Blank time -3 -1		Incubation	Time 0	29
		Reaction		29
· · · ·	Diluent	54	Reagent Vol	
Standard 3µLµLµLµLµL	μι	R1 R2	125 μL	
Increased µL µL	μι	N2	με	
Sample Blank Auto rerun	<u>.</u> .			
Define / Edite Chemestries				
		i	6. I =	
Chem AUML N°			Sample Type	
Chemistry URIC ACID			Print name	
Linearity range (standard) * *		Lin	earity Limit *	
Linearity range (Decreased) * *	Substrate Depletion			
Linearity range (Increased) * *	Mixed Blank Abs *			
R1 Blank Abs * *	On-borad Stability * Jour			
Blank Response * *	Reagent Alarm Limit			
Prozone Check (0) Rtae Check				
Q1 Q2 Q3	Q4	PC		ABS
Calibration setup				
Chem AUML				
Calibration settings	Calibrator Water	Conc. 0	Pos W	Lot No.
Math Model Two-point linear	Cal1	xxx	xxx	
Factor Replicates 2				
Acceptance Limits				
Cal Time Heure				
Slope DiffSD				
Sensitivity Reproductb				
Deter Coeff				
Auto calib.				
□Bottle Changed □Lot Changed □Cal Time				

*Values entered by the operator