

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.

Define / Edite Chemestries

Chem	<input type="text" value="TGML"/>	N°	<input type="text"/>	Sample Type	<input type="text" value="Serum"/>
Chemistry	<input type="text" value="TRIGLYCERIDES"/>	Print name	<input type="text"/>		
Reaction Type	<input type="text" value="Endpoint"/>	Reaction Direction	<input type="text" value="Positive"/>		
Pri Wave	<input type="text" value="505 nm"/>	Sec Wave	<input type="text"/>		
Unit	<input type="text" value="g/dL"/>	Decimal	<input type="text" value="0"/>		
Blank time	<input type="text" value="-3"/>	<input type="text" value="-1"/>	Incubation Time	<input type="text" value="0"/>	
			Reaction Time	<input type="text" value="60"/>	<input type="text" value="62"/>
Sample Vol	<input type="text" value="2"/>	Aspirated	<input type="text"/>	Diluent	<input type="text"/>
Standard	<input type="text"/>	<input type="text"/>	<input type="text"/>	Reagent Vol	
Decreased	<input type="text"/>	<input type="text"/>	<input type="text"/>	R1	<input type="text" value="200"/>
Increased	<input type="text"/>	<input type="text"/>	<input type="text"/>	R2	<input type="text"/>
	<input type="text"/>	<input type="text"/>	<input type="text"/>		
<input type="checkbox"/> Sample Blank		<input type="checkbox"/> Auto rerun			

Define / Edite Chemestries

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Chemistry	<input type="text" value="TRIGLYCERIDES"/>	Print name	<input type="text"/>		

Linearity range (standard)	<input type="text" value="*"/>	<input type="text" value="*"/>	Linearity Limit	<input type="text" value="*"/>
Linearity range (Decreased)	<input type="text" value="*"/>	<input type="text" value="*"/>	Substrate Depletion	<input type="text"/>
Linearity range (Increased)	<input type="text" value="*"/>	<input type="text" value="*"/>	Mixed Blank Abs	<input type="text" value="*"/> <input type="text"/>
R1 Blank Abs	<input type="text" value="*"/>	<input type="text" value="*"/>	On-borad Stability	<input type="text" value="*"/> Jour
Blank Response	<input type="text" value="*"/>	<input type="text" value="*"/>	Reagent Alarm Limit	<input type="text"/>
Twin Chemistry	<input type="text"/>	<input type="checkbox"/> Enzyme Linear Extension		

<input type="checkbox"/> Prozone Check	<input type="checkbox"/> Rtae Check				
Q1 <input type="text"/>	Q2 <input type="text"/>	Q3 <input type="text"/>	Q4 <input type="text"/>	PC <input type="text"/>	ABS <input type="text"/>

Calibration setup

Chem	<input type="text" value="TGML"/>
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Calibration settings	
Math Model	<input type="text" value="Two-point linear"/>
Factor	<input type="text"/>
Replicates	<input type="text" value="2"/>

Acceptance Limits	
Cal Time	<input type="text"/> Heure
Slope Diff	<input type="text"/> SD <input type="text"/>
Sensitivity	<input type="text"/> Reproductb. <input type="text"/>
Deter Coeff	<input type="text"/>

Auto calib.			
<input type="checkbox"/> Bottle Changed	<input type="checkbox"/> Lot Changed	<input type="checkbox"/> Cal Time	

Calibrator	Conc.	Pos	Lot No.
Water	0	W	
Cal1	xxx	xxx	

*Values entered by the operator