Molecular Dimensions

Soluble Proteins

SG1 Screen HT-96 / FX-96 MD1-89 / MD1-89-FX

From the CSRIO - the C3 ShotGun (SG1) Screen - a set of 96 conditions that occur with highest, non-redundant frequency amongst all PDB deposits.

SG1 Screen uses the ShotGun approach to gather all the most successful conditions from all the early commercially available screens.

MD1-89 is presented as 96 x 1 mL conditions/MD1-89-FX is presented as 96 x 100 µL conditions

Features of SG1 (ShotGun) Screen:

- 96 of the most successful conditions from all the early commercially available screens.
- Provides a great start for easy optimization.
- Save money and time.

Introduction

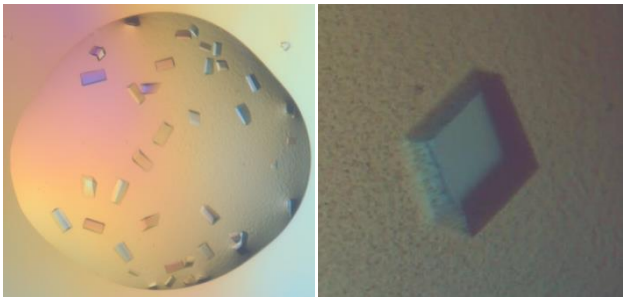
Commercial crystallization screening offers more than 15000 crystallization conditions for screening of new crystallization targets. Some conditions have been far more successful than others.

SG1 (ShotGun) Screen is designed by Janet Newman *et al* from **CSIRO** and represents the most successful, non-redundant frequently reported crystallization conditions* from all the early commercially available crystallization screens.

The term "shotgun screening" was coined early in the Structural Genomics era and refers to the process of setting up experiments using pre-mixed cocktails until a crystal of sufficient quality is obtained. The best place to start screening is within the context of previously successful crystallization space. "*Although only 14% of successful crystallization conditions from, the PDB are identical to a commercial condition, almost 40% of the PDB conditions can be obtained by trivial optimization of a commercial cocktail.*" (Fazio *et al*) So this is a reasonable place to commence screening.

However, that does leave 60% of deposits not covered in this screen and requiring additional screening tools.

Molecular Dimensions has always recommended JCSG-*plus*™ and PACT *premier*™ as powerful non-redundant screens containing 384 conditions that combine a sparse matrix approach with a systematic screen to give maximum information. The recent introduction of Morpheus®, MIDAS™ and The PGA Screen™ provide the opportunity to explore even wider crystallization space with the use of ligand screening and novel precipitants.



Images of HSP90 (N-term domain) crystallized in C3 for MecRx from SG1 (ShotGun) Screen *(courtesy of J.Newman, CSIRO)*

Reference:

Fazio VJ, Peat TS & Newman J (2014). A drunken search in Crystallization Space. Acta Cryst. F70:1303-1311

**These are conditions from commercially available crystallization screens that have been included in the REMARK280 field of the PDB ID code. It is therefore, biased slightly towards the earlier generation of classic screens, and not the later releases, such as Morpheus, MIDAS etc. REMARK280 is a non-mandatory field in the PDB record and should only contain data associated with the crystallization cocktail and not the chemistry associated with protein formulation, the cryoprotectant or soaking solutions.*

Formulation Notes:

SG1 Screen reagents are formulated using ultrapure water (>18.0 MΩ) and are sterile-filtered using 0.22 µm filters. No preservatives are added.
Final pH may vary from that specified on the datasheet. Molecular Dimensions will be happy to discuss the precise formulation of individual reagents.
Individual reagents and stock solutions for optimization are available from Molecular Dimensions.

Enquiries regarding SG1 Screen formulation, interpretation of results or optimization strategies are welcome. Please e-mail, fax or phone your query to Molecular Dimensions.

Contact and product details can be found at www.moleculardimensions.com

Manufacturer's safety data sheets are available from our website or by scanning the QR code here:

Re-Ordering details:

Catalogue Description		Catalogue Code
SG1 Screen	96 x 10 mL	MD1-88
SG1 Screen HT-96	96 x 1 mL	MD1-89
SG1 Screen FX-96	96 x 100 µL	MD1-89-FX
Eco Screens		
SG1 Eco Screen	96 x 10 mL	MD1-88-ECO
SG1 Screen HT-96 Eco Screen	96 x 1 mL	MD1-89-ECO
Single Reagents		
SG1 Screen single reagent	100 mL	MDSR-88-tube number
SG1 Screen HT-96 single reagent	100 mL	MDSR-89-well number

For SG1 Screen stock reagents visit our Optimization page on our website.

*Developed by Janet Newman (CSIRO) and manufactured under licence by Molecular Dimensions Ltd. Molecular Dimensions operates an ethical approach to all its products by making sure the inventors of its products receive the appropriate acknowledgments/rewards for their hard work. We hope you appreciate their hard work too and only buy the 'real-thing' anything else that is not licenced does not acknowledge or support the inventor and institute.

Molecular Dimensions

Soluble Proteins

SG1 Screen HT-96 Conditions A1 - D12 MD1-89 / MD1-89-FX

Well #	Conc.	Salt1	Conc.	Salt2	Conc.	Buffer	pH	Conc.	Precipitant1	Conc.	Precipitant2
A1	0.2 M	Magnesium chloride hexahydrate			0.1 M	Tris	8.5	30 % w/v	PEG 4000		
A2	2.0 M	Ammonium sulfate									
A3	0.2 M	Sodium acetate trihydrate						20 % w/v	PEG 3350		
A4	2.0 M	Ammonium sulfate			0.1 M	Tris	8.5				
A5	0.2 M	Sodium citrate tribasic dihydrate						20 % w/v	PEG 3350		
A6					0.1 M	Sodium HEPES	7.5	20 % w/v	PEG 4000	10 % v/v	2-Propanol
A7	2.0 M	Ammonium sulfate			0.1 M	Sodium HEPES	7.5	2 % v/v	PEG 400		
A8	1.4 M	Sodium citrate tribasic dihydrate			0.1 M	Sodium HEPES	7.5				
A9	0.2 M	Sodium acetate trihydrate			0.1 M	Tris	8.5	30 % w/v	PEG 4000		
A10	0.2 M	Lithium sulfate			0.1 M	Tris	8.5	30 % w/v	PEG 4000		
A11	4.0 M	Sodium formate									
A12	0.2 M	Magnesium acetate tetrahydrate			0.1 M	Sodium cacodylate	6.5	20 % w/v	PEG 8000		
B1					0.1 M	Bis-Tris	5.5	25 % w/v	PEG 3350		
B2					0.1 M	MES	6.5	12 % w/v	PEG 20000		
B3	0.2 M	Magnesium chloride hexahydrate			0.1 M	Bis-Tris	5.5	25 % w/v	PEG 3350		
B4	0.2 M	Ammonium sulfate			0.1 M	MES	6.5	30 % w/v	PEG 5000 MME		
B5	0.2 M	Calcium chloride dihydrate						20 % w/v	PEG 3350		
B6					0.1 M	Sodium HEPES	7.5	20 % w/v	PEG 10000		
B7	0.2 M	Sodium formate						20 % w/v	PEG 3350		
B8	0.2 M	Ammonium sulfate			0.1 M	Bis-Tris	5.5	25 % w/v	PEG 3350		
B9	1.6 M	Sodium citrate tribasic dihydrate									
B10	0.2 M	Calcium chloride dihydrate			0.1 M	Sodium HEPES	7.5	28 % v/v	PEG 400		
B11	0.2 M	Ammonium chloride						20 % w/v	PEG 3350		
B12	0.2 M	Magnesium formate dihydrate						20 % w/v	PEG 3350		
C1	0.2 M	Ammonium sulfate			0.1 M	Sodium acetate	4.6	25 % w/v	PEG 4000		
C2	1.4 M	Sodium malonate dibasic monohydrate pH 7.0									
C3	0.2 M	Lithium sulfate			0.1 M	Bis-Tris	5.5	25 % w/v	PEG 3350		
C4	0.2 M	Potassium sodium tartrate tetrahydrate						20 % w/v	PEG 3350		
C5	0.2 M	Ammonium sulfate			0.1 M	Sodium cacodylate	6.5	30 % w/v	PEG 8000		
C6	2.0 M	Ammonium sulfate			0.1 M	Sodium acetate	4.6				
C7					0.1 M	Sodium HEPES	7.5	25 % w/v	PEG 3350		
C8	0.2 M	Magnesium chloride hexahydrate			0.1 M	Bis-Tris	6.5	25 % w/v	PEG 3350		
C9	0.2 M	Magnesium chloride hexahydrate			0.1 M	Tris	8.5	25 % w/v	PEG 3350		
C10	0.2 M	Magnesium chloride hexahydrate			0.1 M	Sodium HEPES	7.5	25 % w/v	PEG 3350		
C11	0.2 M	Sodium acetate trihydrate			0.1 M	Sodium cacodylate	6.5	30 % w/v	PEG 8000		
C12	0.2 M	Sodium acetate trihydrate			0.1 M	Bis-Tris	5.5	25 % w/v	PEG 3350		
D1	1.5 M	Lithium sulfate			0.1 M	Sodium HEPES	7.5				
D2					0.1 M	Sodium citrate	5.5	20 % w/v	PEG 3000		
D3								25 % w/v	PEG 1500		
D4	0.2 M	Potassium thiocyanate						20 % w/v	PEG 3350		
D5	0.2 M	Sodium acetate trihydrate			0.1 M	Sodium cacodylate	6.5	18 % w/v	PEG 8000		
D6	0.2 M	Lithium sulfate			0.1 M	Sodium HEPES	7.5	25 % w/v	PEG 3350		
D7	0.2 M	Ammonium sulfate						30 % w/v	PEG 8000		
D8					0.1 M	Bis-Tris	6.5	20 % w/v	PEG 5000 MME		
D9	0.2 M	Ammonium sulfate			0.1 M	Sodium acetate	4.6	30 % w/v	PEG 2000 MME		
D10	0.2 M	Lithium sulfate			0.1 M	Bis-Tris	6.5	25 % w/v	PEG 3350		
D11					0.1 M	Sodium acetate	4.6	8 % w/v	PEG 4000		
D12	2.0 M	Ammonium sulfate			0.1 M	Bis-Tris	6.5				

Molecular Dimensions

Soluble Proteins

SG1 Screen HT-96 Conditions E1 - H12 MD1-89 / MD1-89-FX

Well #	Conc.	Salt1	Conc.	Salt2	Conc.	Buffer	pH	Conc.	Precipitant1
E1	2.0 M	Ammonium sulfate			0.1 M	Bis-Tris	5.5		
E2								25 % w/v	PEG 3350
E3	0.2 M	Magnesium chloride hexahydrate			0.1 M	Sodium HEPES	7.5	30 % v/v	PEG 400
E4	2.0 M	Ammonium sulfate			0.1 M	Sodium HEPES	7.5		
E5	3.5 M	Sodium formate							
E6	1.6 M	Magnesium sulfate heptahydrate			0.1 M	MES	6.5		
E7	0.2 M	Magnesium chloride hexahydrate						20 % w/v	PEG 3350
E8	0.2 M	Ammonium sulfate						30 % w/v	PEG 4000
E9	0.1 M	Potassium thiocyanate						30 % w/v	PEG 2000 MME
E10	0.2 M	Sodium malonate dibasic monohydrate pH7						20 % w/v	PEG 3350
E11	2.0 M	Sodium formate			0.1 M	Sodium acetate	4.6		
E12	0.2 M	Ammonium sulfate			0.1 M	Sodium HEPES	7.5	25 % w/v	PEG 3350
F1	0.2 M	Potassium sodium tartrate tetrahydrate	2.0 M	Ammonium sulfate	0.1 M	Sodium citrate	5.6		
F2	0.2 M	Sodium acetate trihydrate			0.1 M	Sodium HEPES	7.5	25 % w/v	PEG 3350
F3	0.2 M	Ammonium sulfate						20 % w/v	PEG 3350
F4	1.0 M	Sodium citrate tribasic dihydrate			0.1 M	Sodium cacodylate	6.5		
F5	0.2 M	Ammonium sulfate			0.1 M	Bis-Tris	6.5	25 % w/v	PEG 3350
F6	0.2 M	Ammonium nitrate						20 % w/v	PEG 3350
F7	0.2 M	Sodium thiocyanate						20 % w/v	PEG 3350
F8	0.2 M	Potassium nitrate						20 % w/v	PEG 3350
F9					0.1 M	Sodium HEPES	7.5	20 % w/v	PEG 8000
F10	0.2 M	Magnesium acetate tetrahydrate						20 % w/v	PEG 3350
F11					0.1 M	Bis-Tris	6.5	25 % w/v	PEG 3350
F12	0.02 M	Calcium chloride dihydrate			0.1 M	Sodium acetate	4.6	30 % v/v	MPD
G1	0.2 M	Sodium acetate trihydrate			0.1 M	MES	6.0	20 % w/v	PEG 8000
G2	0.2 M	Sodium sulfate						20 % w/v	PEG 3350
G3	0.01 M	Zinc sulfate heptahydrate			0.1 M	MES	6.5	25 % v/v	PEG 550 MME
G4	0.2 M	Sodium tartrate dibasic dihydrate						20 % w/v	PEG 3350
G5								60 % v/v	T-mate pH 7.0
G6	0.5 M	Ammonium sulfate	1.0 M	Lithium sulfate	0.1 M	Sodium citrate	5.6		
G7								30 % w/v	PEG 1500
G8	0.2 M	Magnesium chloride hexahydrate			0.1 M	Tris	8.5	20 % w/v	PEG 8000
G9	0.2 M	Ammonium tartrate dibasic						20 % w/v	PEG 3350
G10	0.2 M	Sodium fluoride						20 % w/v	PEG 3350
G11	0.2 M	Sodium chloride	2.0 M	Ammonium sulfate	0.1 M	Sodium cacodylate	6.5		
G12	0.1 M	Sodium chloride	1.6 M	Ammonium sulfate	0.1 M	Sodium HEPES	7.5		
H1	0.2 M	Ammonium formate						20 % w/v	PEG 3350
H2	0.2 M	Lithium citrate tribasic tetrahydrate						20 % w/v	PEG 3350
H3	0.2 M	Ammonium iodide						20 % w/v	PEG 3350
H4	0.2 M	Sodium acetate trihydrate			0.1 M	Bis-Tris	6.5	25 % w/v	PEG 3350
H5								30 % w/v	PEG 4000
H6					0.1 M	Tris	8.5	25 % w/v	PEG 3350
H7	0.2 M	Ammonium fluoride						20 % w/v	PEG 3350
H8	0.1 M	Sodium acetate trihydrate			0.1 M	Bis-Tris	5.5	17 % w/v	PEG 10000
H9	0.2 M	Sodium acetate trihydrate			0.1 M	Imidazole	8.0	10 % w/v	PEG 8000
H10	0.2 M	Ammonium sulfate			0.1 M	Tris	8.5	25 % w/v	PEG 3350
H11					0.1 M	CHES	9.0	20 % w/v	PEG 8000
H12	4.3 M	Sodium chloride			0.1 M	Sodium HEPES	7.5		

Abbreviations: Bis-Tris; Bis-(2-hydroxyethyl)imino-tris(hydroxymethyl)methane, **CAPS**; N-Cyclohexyl-3-aminopropanesulfonic acid**, CHES**; 2-(N-Cyclohexylamino)ethane Sulfonic Acid**, HEPES;** 2-(4-(2-Hydroxyethyl)-1-piperazinyl)ethanesulfonic Acid, **Sodium HEPES**; 2-(4-(2-Hydroxyethyl)-1-piperazinyl)ethanesulfonic Acid Sodium Salt, **MES**; 2-(N-morpholino)ethanesulfonic acid, **MPD;** 2,4-methyl pentanediol, **PEG**; Polyethylene glycol, **T-mate;** Sodium malonate dibasic monohydrate, Ammonium citrate tribasic, Succinic acid, DL-Malic acid, Sodium acetate trihydrate, Sodium formate, Ammonium tartrate dibasic, **Tris**; 2-Amino-2-(hydroxymethyl)propane-1,3-diol.