



2228.HK

XtalPi Holdings

H1 2026 Investor Presentation

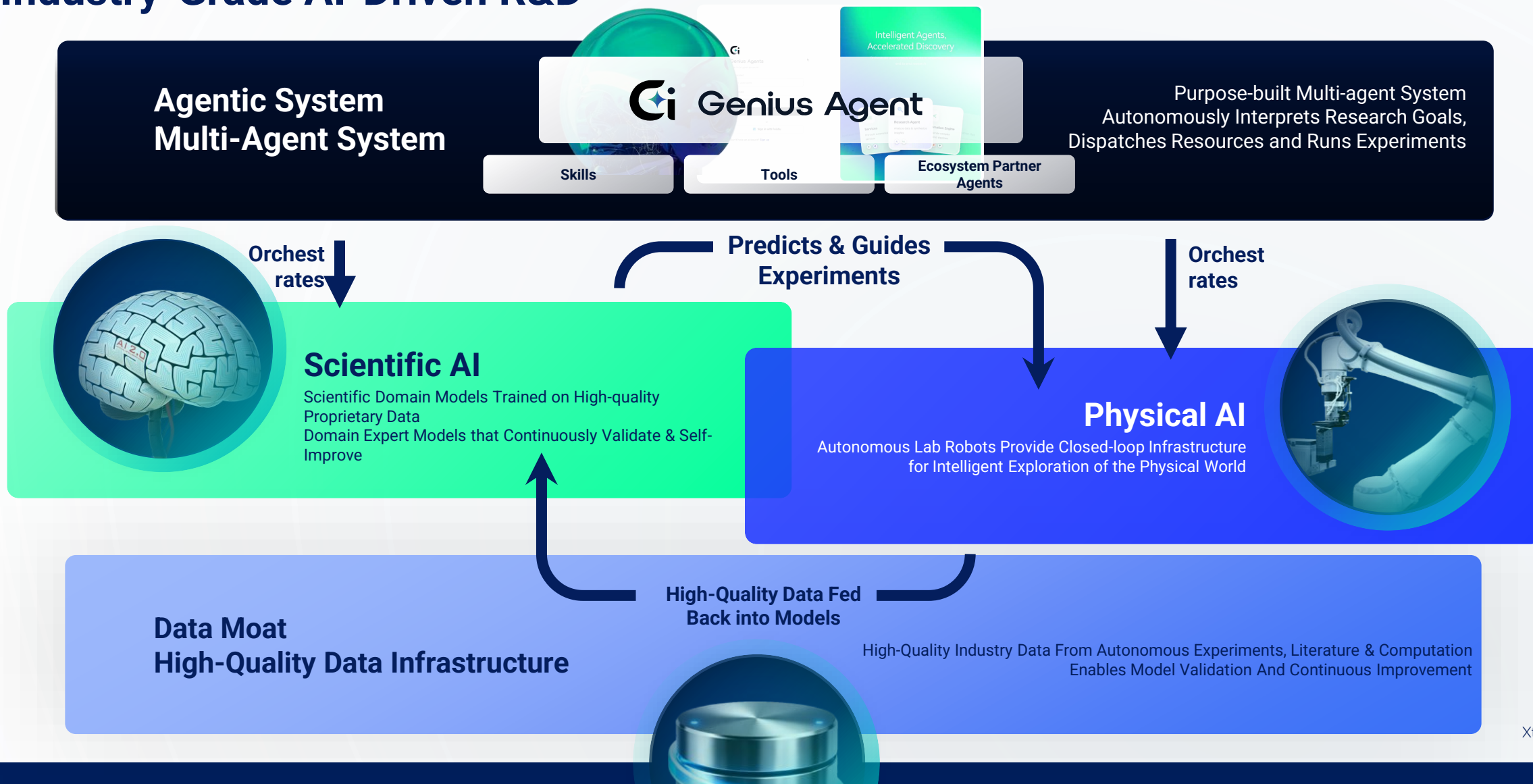


AI4S: Connecting Digital & Physical Worlds to Enable Autonomous Scientific Discovery

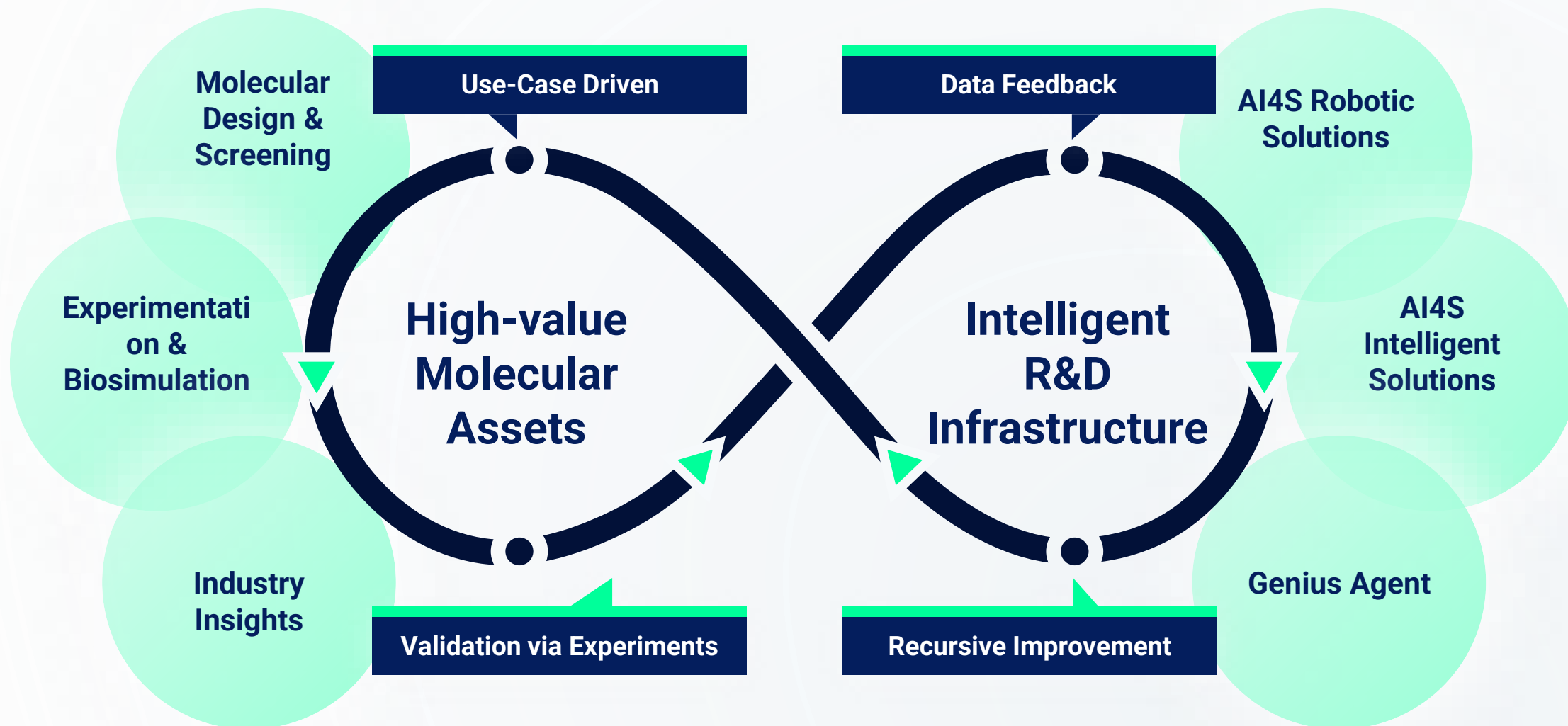
XtalPi's Current Capabilities



Technology Engine: A Four-Layer Core Architecture Underpinning Industry-Grade AI-Driven R&D



Two Businesses, One Self-Improving Closed Loop



Business Highlights: Multiple Major Breakthroughs Across the Group

Leading the Development of AI4S Infrastructure

AI4S Intelligent Solutions
+136.4% YoY

AI Drug R&D Platform with the Broadest Modality Coverage

Small molecules, large molecules, peptides and oligonucleotides
Four core tech platforms

Rapid Build-out of Differentiated, Extensive Pipeline

Partnered and proprietary portfolio
Over 20 programs at PCC

"Real-world" AI Breaking Through Conventional Drug R&D Bottlenecks

Continued validation of
complex, problem-solving capabilities

Strong Recognition and Comprehensive Collaboration with Leading Customers

Covering
platform services, model licensing, pipeline deals and AI4S infrastructure deployment

Launch of a Self-Evolving AI Retrosynthesis Chemistry System

"Chemical hallucination" rate
1/6 that of leading large models

First Comprehensive Open Scientific Research Platform to Connect Physical AI

XtalPi Science
autonomous scientific discovery platform

Deploying Leading Biological Models

Virtual cells, organ-on-chip and organoids
boost the translational success rate of preclinical R&D

Outlook: From Expanding Orders to R&D Transformation, Full Industry-Chain Presence Opens Long-Term Growth Space

Current

Order Expansion

- AI-driven drug discovery is driving demand for **new molecule synthesis and R&D data**
- AI-native experimentation systems and intelligent laboratories winning **orders from leading pharmaceutical companies**

Medium Term

Platform + Assets

- Pharmaceutical companies in China and overseas need to replenish pipelines, lifting demand for advanced R&D services and high-quality assets
- **Platform technology services** generate stable revenue; **asset deals** monetizes high-quality proprietary pipelines

Mid-to-Long Term

Dual Engine

- Leverage the efficiency of AI-driven R&D to continue incubating high-quality proprietary pipelines
- As pipelines move into regulatory filings and clinical development, **R&D services and proprietary new drugs form a dual-engine** that takes growth beyond the limits of a services-only model.

Long Term

A New R&D Model

- AI4S is driving pharmaceutical R&D from **experience-based trial and error toward a data- and intelligence-driven model**
- End-to-end smart R&D reduces costs and improves efficiency; **proprietary data** and **algorithmic advantages** build a durable moat

2026 H1 Financial Key Numbers

RMB mn	2026 H1	2025 H1	YoY
Revenue	394	517	(24%)
Revenue (Excl. Major Pipeline BD)	264	152	+74%
R&D Expenses	368	222	+66%
(Loss) / Profit for the Period	(225)	76	N/A
Adjusted Net (Loss) / Profit	(106)	142	N/A

Cash Balance*: As of June 30, 2026, the Company's total cash balance amounted to RMB 8,671.2 million.

*Cash balance includes our cash and cash equivalents, bank deposits, the current portion of financial assets measured at FVTPL and restricted cash as of 30 June 2026

Excluding Major Pipeline BD Deals: Organic Revenue Grew Substantially, Driving a Significant Narrowing of the Loss Margin

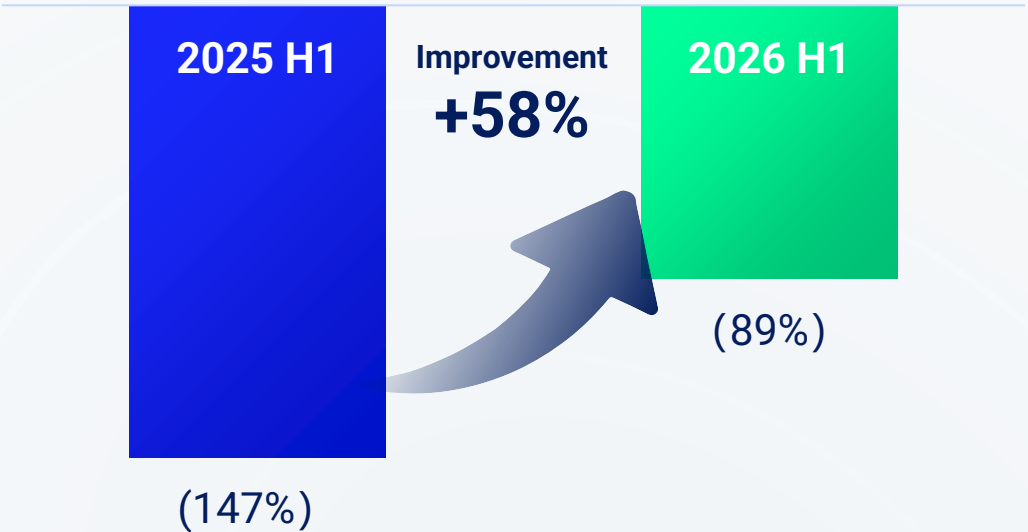
Organic Revenue Grew Substantially

RMB mn



Loss Ratio² Narrowed Substantially

RMB mn



RMB mn

Adjusted net profit / (loss)
Adjusted net profit / (loss) (excluding major pipeline BD)

	2025 H1	2026 H1
Adjusted net profit / (loss)	142	(106)
Adjusted net profit / (loss) (excluding major pipeline BD)	(224)	(235)

¹Year-on-year revenue growth rate excluding major pipeline BD

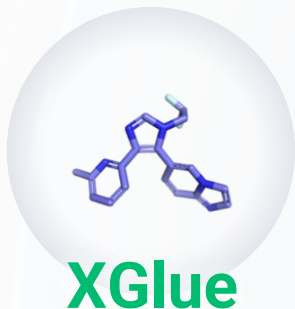
²Loss ratio = Non-IFRS net loss excluding major pipeline BD ÷ revenue excluding major pipeline BD



Drug Discovery Solutions

Four Modality Platforms: Proprietary Data + Generative/Predictive Models Building Differentiated Drug-Discovery Capability

Small Molecule



- Achieved a molecular glue virtual compound library covering millions of molecules, along with a physical scaffold library of tens of thousands of compounds
- Enables synthesis and validation of hundreds to thousands of compounds per week

Large Molecule



- Three engines: XtalFold® structure modeling, XenProT® generative AI, Xentient® discriminative AI
- AtlaX data foundation has advantage over public datasets on key data types

Peptides



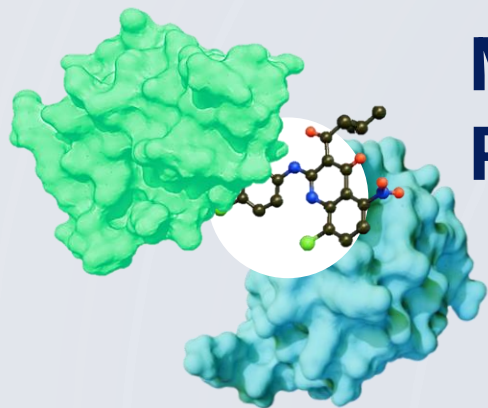
- Core models deliver industry-leading performance in cyclic peptide membrane permeability assessment and generation tasks
- A database of more than 5,000 non-natural amino acids has been established

Oligonucleotides



- Proprietary siFormer architecture jointly designs sequence and chemical modifications, significantly improving generalization to unknown targets
- Supports global optimization of dual-target siRNA in a single molecule; compatible with small-molecule, LNP, antibody and peptide delivery; supports hepatic and extrahepatic targeting

Small Molecules: Million-Scale Library Used to Build Molecular Glue Platform; Multiple Autoimmune Pipelines Achieving Picomolar-level Breakthrough



Molecular Glue Platform

2.58M

Virtual
compound
library

40+

CRBN-binding
scaffolds

100K+

In-stock
molecular
building blocks

Autoimmune-focused Picomolar-level Breakthroughs in Select Molecules

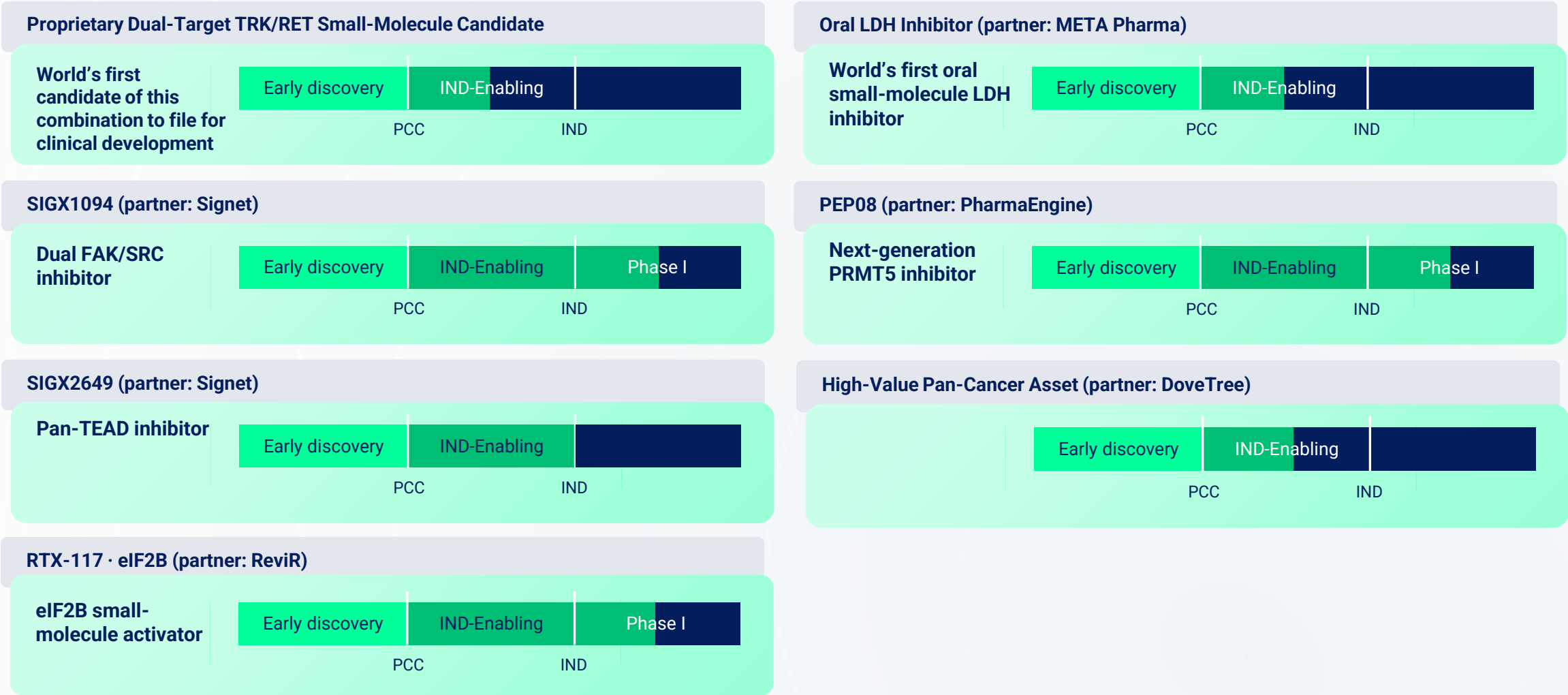
pM

Optimized to picomolar degradation
activity within one quarter

**Multi-
target**

Hit compounds identified against
multiple autoimmune targets

Small-Molecule Pipeline: Core Programs in Clinical Development, Multiple Potential FIC/BIC Assets Entering IND Stage



Biologics (Ailux): World-Leading AI-Native Antibody Platform

XtalFold™ Structure Modeling



Prediction of antigen-antibody complexes and other structures

Integrated with generation and discriminative modules

Supports epitope identification and candidate triage

XenProT™ Generative AI



Molecular Generation and Design

Covers affinity maturation, pH-sensitive engineering and bispecific antibody design

Xentient™ Discriminative AI



Closed-Loop Functional Evaluation and Optimization

Works with structure modeling and generation

Supports candidate selection, multi-objective optimization and other applications

AtlaX™ Data Foundation

Proprietary Wet-Lab System

High-throughput Data Generation

LuxSight™ Patent-Mining Agent

Across affinity, complex structures, antigen-antibody pairing and native heavy/light-chain sequences:

Several-fold to tens-fold scale advantage on each data type

Proprietary assays generate functional labels rarely covered by public data, including polyreactivity, stability and immunogenicity:

Provides a differentiated data moat for complex antibody development

Biologics (Ailux) Pipeline: Clear Differentiation, Preclinical Development Advancing Steadily

ALX001

TL1A×IL-23: Bispecific

Biological Mechanism & Opportunity

- Raises the efficacy ceiling: dual blockade can synergistically suppress inflammation
- Beyond Th17: TL1A blockade may also ease tissue fibrosis

Differentiation

- Anti-TL1A mAb: unique MoA, low immunogenicity risk; does not block TL1A binding to DcR3
- Bispecific: potent anti-inflammatory activity
- Potential BIC, targeting Ph. I in 2027

Fc-silenced
Half-life extended

Anti-IL-23p19 IgG

Anti-TL1A scFvs

Indications



ALX002

CD19×BCMA×CD3: Trispecific TCE

Biological Mechanism & Opportunity

- Clinically validated MoA: immune reset and durable remission off treatment
- More convenient and scalable than CAR-T

Differentiation

- Deep, potent B-cell depletion with only moderate cytokine release
- Proprietary TCE format with a masked anti-CD3 arm for improved safety and developability
- All three binding arms are fully proprietary
- Potential BIC, targeting Ph. I in 2027

Anti-BCMA

Fc-silenced

Anti-CD19

Indications



ALX005

Long-acting FcRn blocker for autoantibody-mediated disease

Biological Mechanism & Opportunity

- AI-driven epitope focusing and XtalFold-guided triage deliver a long-acting FcRn blocker with exceptionally low safety risk
- Viscosity prediction supports convenient SC dosing across autoantibody-mediated disease

Differentiation

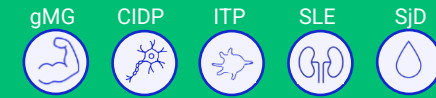
- Unique bispecific design: anti-FcRn IgG × anti-HSA VHH
- Exceptionally low safety risk and suitable for subcutaneous administration
- Potential BIC; Phase I targeted for 2027

Fc-silenced

Anti-FcRn IgG

Anti-HSA VHH

Indications



Peptides: New Product Achieves GRAS Status; Brain-Delivery and Oral Cyclic Peptides Target PCC

HELM-DIFF and related models

Unlocking the NCAA design space

Proprietary NCAA library **5,000+**

HELM-DIFF membrane permeability: **SOTA**

All-Atom Dynamic 3D, RMSD **0.76 Å**

Brain-delivery / oral cyclic-peptide pipeline

PCC targeted for 2027

Brain delivery

In vivo optimization underway

PCC expected
1H27

Oral cyclic peptide for autoimmunity

Hit in 2 months on a selected target; now in Hit-to-Lead

PCC expected mid-2027

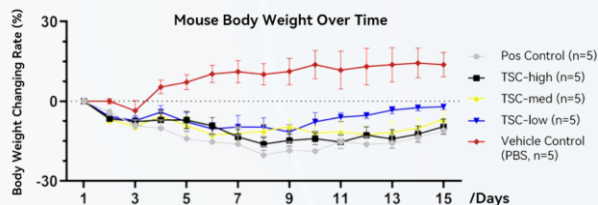
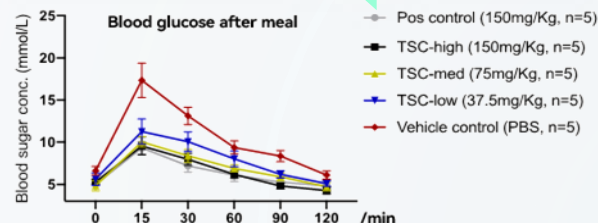
Tensotide™

US Self-Affirmed GRAS status
Enabling use in food and dietary supplements

Strong animal-model data

In mice, within two weeks

~30% blood glucose reduction; 5-10% weight loss



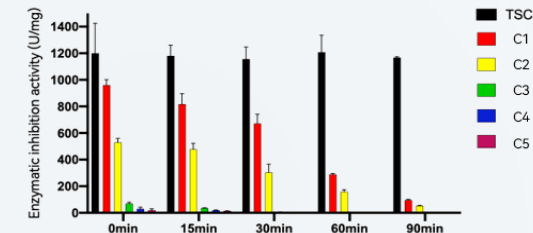
High in vivo stability

At 40°C,

Tensotide™ highly stable across extreme pH and multiple digestive enzymes

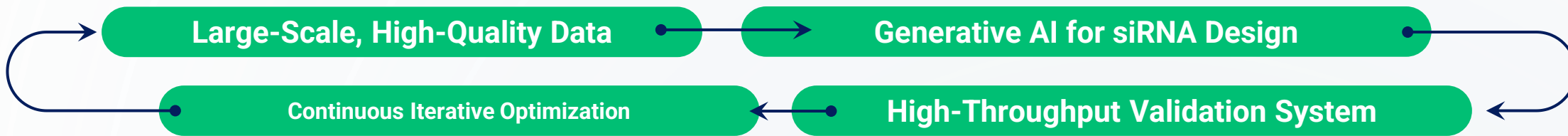
4°C固体	99.72%
40 °C水溶液	99.58%
PH=13(1H)	99.50%
PH=1(1H)	99.72%
INITIATION	99.83%

Comparison of TSC with 5 other commercial products



Oligonucleotides: Platform R&D Efficiency Up >2X; IgA Nephropathy Program Delivered Strong NHP Efficacy Data in 7 Months

Kodexia™ siRNA De Novo Design Platform



Platform Capabilities

AI-Native Design

Deeply integrates first-principles biological mechanism modeling with GenAI and, supported by high-throughput automation, builds a closed-loop iteration between dry and wet experiments.

Tens of Thousands of Wet-Lab Data Points

Covers multiple cell lines, multiple targets, and multi-level, multimodal data from in vitro to in vivo; and has formed a diverse siRNA chemical modification database.

siFormer: Joint Sequence × Modification Design

Generates candidates while recommending personalized modifications; supports global optimization of single-molecule dual-target siRNA; compatible with small-molecule, LNP, antibody and peptide delivery, covering hepatic and extrahepatic tissues.

Efficiency & Hit Rate

≥2X

Using publicly comparable metric, platform R&D efficiency is more than 2X that of traditional methods.

+266%

Molecular property prediction accuracy improved by approximately 266%, establishing predictive capabilities related to in vivo efficacy.

>50%

Across multiple pipelines, >50% of first-round designed molecules showed stronger in vivo activity than the positive control.

Pipeline Portfolio

6

6 pipelines established; over half have completed in vivo efficacy evaluation

Fastest program has entered PCC

IgA Nephropathy Lead Program

Lead program obtained NHP efficacy data about **7 months** after launch, with activity and durability better than a same-target clinical-stage reference

Metabolism

Single-target and dual-target weight-loss siRNA programs are being advanced in parallel

Pipeline Progress Overview

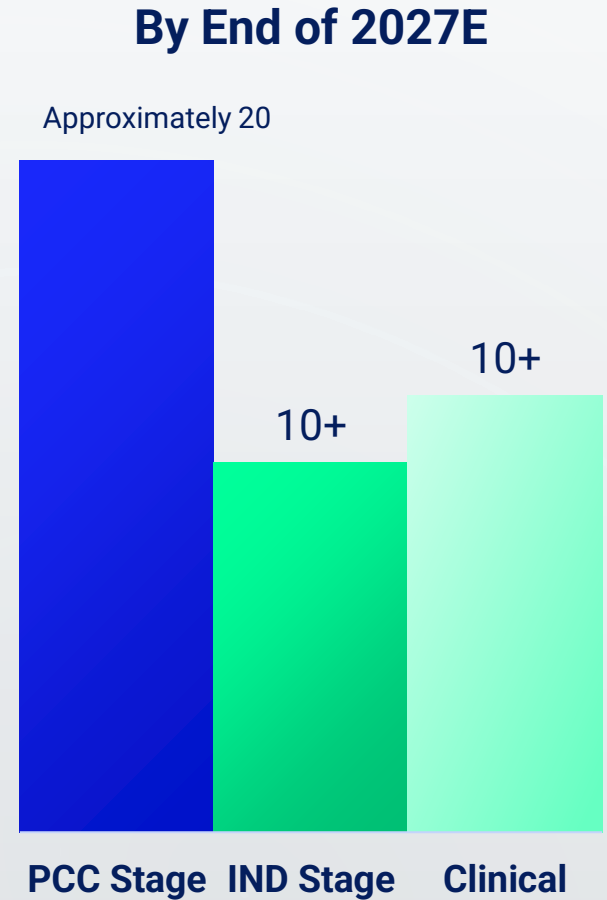
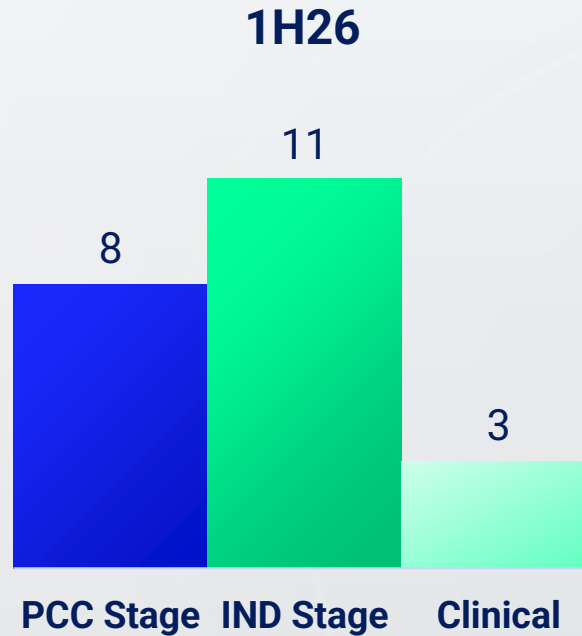
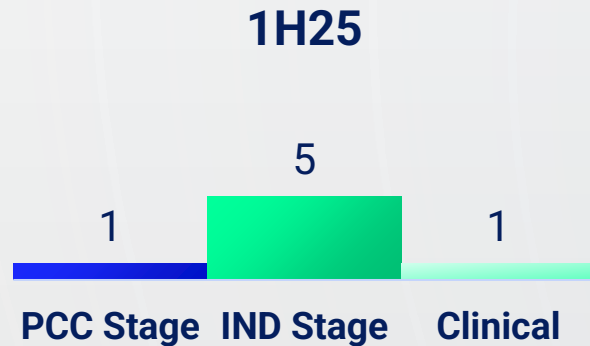
 Self developed

 Co-developed

Modality	Pipeline	Target/MoA	Indication	PCC	IND	Ph. I	Ph. II/III	Commercialization	Rights / partner	Next milestones	Remarks
Bispecific Ab	ALX001	TL1A/IL-23p19	Autoimmune disease						Self-developed	Ph. I in 2027	Potential BIC
Trispecific Ab	ALX002	CD19/BCMA/CD3	Autoimmune disease						Self-developed	Ph. I in 2027	Potential BIC
Bispecific Ab	ALX005	FcRn/HSA	Autoimmune disease						Self-developed	Ph. I in 2027	Potential BIC
Small molecule	XTN004	TRK/RET	Intestinal pain						Self-developed	China/US IND filings in 2H26	—
Small molecule	SIGX1094	FAK/SRC	Diffuse gastric cancer (DGC)						Signet	Ph. II in 2027	Potential FIC, Prix Gallien nominee, ODD, Fast Track
Small molecule	RTX-117	eIF2B	Charcot-Marie-Tooth disease (CMT); vanishing white matter disease (VWM)						ReviR	Ph. II in 2027	Potential FIC, ODD
Small molecule	PEP08	PRMT5	Solid tumour						PharmaEngine		Potential BIC
Small molecule	SIGX2649	TEAD	Solid tumours (mesothelioma, liver cancer, lung cancer, etc.)						Signet	Ph. I in 2H26	Potential FIC/BIC, key data presented at 2026 AACR
Small molecule	MP-5342	LDH	Inflammatory bowel disease (IBD)						META	Ph. I in 2027	Potential FIC
Small molecule	-	-	Oncology						DoveTree	IND filing in 2027	—
Cell	META10-19	CD19, IL-10	Relapsed/refractory CD19+ B-cell hematologic malignancy						Leman	Ph. I in 2H26	Potential BIC
Cell	META10-19	CD19, IL-10	Refractory B-cell autoimmune disease						Leman	Ph. I in 2H26	Potential BIC
Cell	IL-10 GPC3 CAR-T	GPC3, IL-10	GPC3+ liver cancer						Leman	First solid-tumour clinical data readout in 2H26	Potential BIC
Cell	IL-10 EGFRvIII CAR-T	EGFRvIII, IL-10	EGFRvIII+ glioma						Leman	First solid-tumour clinical data readout in 2H26	Potential BIC
Cell	IL-10 Claudin18.2 CAR-T	Claudin18.2, IL-10	Claudin18.2+ gastric cancer						Leman	IIT Beginning in 2H26	Potential BIC
Peptide	Px-002C	Single target	Glucose-lowering	 US Self-Affirmed GRAS					Self-developed		Consumer product
Peptide	Px-003C	Single target	Lipid-lowering						Self-developed		Consumer product
Peptide	Px-001C	Single target	Cosmetics						Self-developed	New cosmetic ingredient filing in 2H27	Consumer
Oligonucleotide	XS01	Single target	IgA nephropathy / PNH						Self-developed		Potential BIC

AI-Native R&D Systems Empowering Progress; Multimodal Assets are Entering a Period of Explosive Value Realization

Number of in-house and partner-enabled pipelines



Landing Major Partnerships in Rapid Succession, Diversified Profit Model Continues to Progress

Reached an AI drug discovery strategic collaboration with an internationally renowned biopharmaceutical company, totaling over USD 400 million



维昇药业
VISEN

Key cooperation worth tens of millions of RMB with Visen Pharmaceuticals



Signed a major strategic cooperation agreement with Sunshine Lake Pharma



DoveTree

The collaboration with DoveTree is progressing well, the Group has received the second payment as stipulated under the final agreement.

晶泰科技收到DoveTree第二笔付款，首个肿瘤项目推进至IND-enabling研究阶段

XtalPi
晶泰科技

Signed a global strategic cooperation and platform licensing agreement with Gan & Lee Pharmaceuticals for AI-driven innovative peptide drug R&D

甘李
GL



Biological Simulation Platforms: Shortening the Validation Pathway from Molecules to Tissues

Moving human-relevant efficacy and safety assessment earlier
Improving candidate translatability

Biosimulation based on new approach methodologies (NAMs) to complement
and partially replace experiments with low translational relevance

Target

Hit

Lead

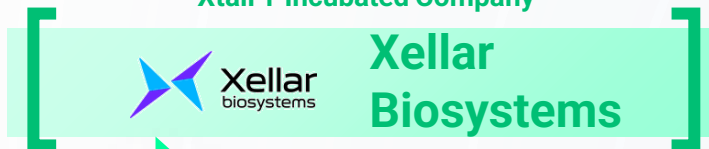
PCC

Preclinical

Clinical

Organ-on-a-Chip

XtalPi-Incubated Company



Delivered and implemented projects for Sanofi / Pfizer;
completed a Series A round of approximately RMB 400 million

Self-Developed Automated Organ-on-a-Chip Production Line x Vertical AI Foundation Model for Biology

Accumulating human-derived 3D wet-lab data;
Building a human biology atlas

Commercialization & Industry Collaboration Outcomes

Delivered iDEA-TECH to Sanofi, fully collected the final payment;
completed a key milestone delivery for jointly-developed AI
toxicity prediction system with Pfizer

Regulatory Compliance

Deeply involved in the CDE NAMs multi-center collaborative
validation program and selected for the "Pioneer Program" while
advancing FDA ISTAND model qualification.

Virtual Cells

XtalPi-Incubated Company



AIVC achieves SOTA on three major perturbation
benchmarks

AI Virtual Cell Model OCOO-T

Achieves industry-leading performance on chemical
drugs, gene and cytokine benchmarks

Scientific Discovery Engine: The Popper Project (TPP)

Internal beta applications opening soon

Completed Angel Round Of Tens of Millions of RMB

TPP to begin internal beta testing soon., with
participation from Shunwei Capital, Sequoia China
and Songhe Capital.

Organoids

XtalPi Partner



Over 15 organoid models; high accuracy in toxicity
prediction models

Organoid + AI Platform

More than 15 standardized systems for gene-edited tumor
organoid models, including diffuse gastric cancer, liver
cancer, and lung cancer, for drug development.

Normal Organoids + AI, including Heart / Kidney / Liver

The drug-induced cardiotoxicity prediction model based on cardiac
organoids achieved an accuracy of 81.25%, far exceeding the 43.75% of
traditional methods; the platform is also being gradually expanded to non-
oncology therapeutic pipelines.

Participation in Industry Standards Development

Deeply involved in national organoid standard-setting projects led by the
CDE and the "Innovative Methods Evaluation Center" of the National
Institutes for Food and Drug Control.



AI4S Intelligent Solutions

AI4S Intelligent Robotic Laboratories (Physical AI): Dense Delivery of Flagship Projects; Multiple RMB10-Million-Level Projects in China, Strong Recognition from Leading Overseas Customers

2026

Signed a RMB 10s-of-millions compound storage and management system project with Eli Lilly



Signed a RMB 10s-of-millions HTE platform cooperation project with Eli Lilly



Completed delivery for JW Pharma of a high-throughput automated synthesis workstation, an AI reaction-condition optimization system, and an intelligent analytics platform



Signed a RMB million-level intelligent high-throughput preparation project for mesoporous materials with Fudan University



Collaborating with Peking University and the Dalian Institute of Chemical Physics for high-throughput electrolyte preparation and testing



2025 and Earlier

Pharmaceuticals

Globally-leading Pharmaceutical Company Headquartered in Indianapolis

Intelligent compound management system for micro-powder dispensing

Roche

Intelligent compound storage system

Haleon

Intelligent sample pre-treatment system

Pharmaceutical Company

Automated compound library-building robot

Petrochemicals/Chemicals

Mainland Chinese SOE

Intelligent high-throughput platform for catalytic materials R&D

BASF

Intelligent workstation for formulation stability testing

New Energy

Domestic University Laboratory

Electrolyte R&D automated workstation

Research Institutions

Liangzhu Laboratory

Intelligent biomaterials R&D platform

Fudan University

Organic synthesis methodology automated workstation

Traditional Chinese Medicine

Guangdong Hengqin Laboratory

Traditional Chinese Medicine extraction and separation automated workstation cluster

AI4S Intelligent Services: VAST™ Virtual Compound Library. Helping Customers Expand Chemical Space and Shorten R&D Cycles

Virtual · Automated · Swift · Tailored | Navigating actionable, underexplored chemical space with automation and proprietary data

Product Positioning & Commercial Breakthrough

VAST™ delivers novel, synthesizable compounds with fast turnaround and high cost-effectiveness. **The new VAST Virtual Compound Library business has achieved a breakthrough, securing orders for over 20,000 molecules and opening up an entirely new growth space for the group.** Empowering customers to expand chemical space and shorten R&D cycles.

5 billion+

Total Unique
Compounds

85%+

Synthesis Success
Rate

1–2

Typical Number of
Synthetic Steps

2–4 Weeks

Average Delivery
Time

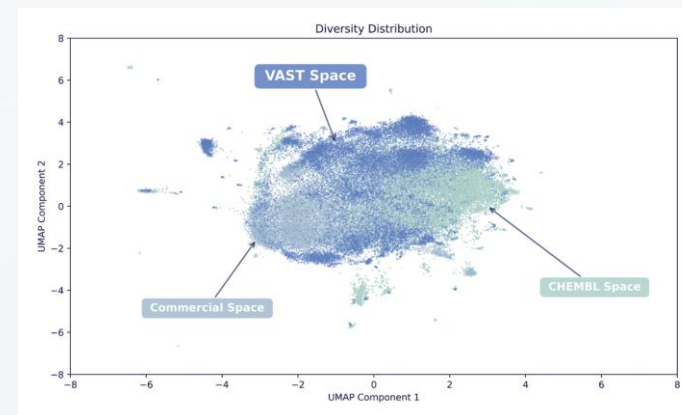
10,000+

New Molecular Building
Blocks

300+

Automated
Workstations

Novel Chemical Space · Diversity Distribution



How the library is built:

- 1–2 step enumeration: 58,000 building blocks × 42 validated reaction protocols
- AI feasibility scoring (proprietary data) filters for synthesizability
- Property filters / novelty checks / PAINS filters
- Diversity subsets and Ro5 / Beyond-Ro5 options; 500 million downloadable
- Automated platforms (dispensing / reaction / analysis / purification, etc.) ensure fast and reliable delivery

AI4S Intelligent Services: Agentic Synthesis Boosts Throughput by Orders of Magnitude, Reaction Success Rate >90%



Scientific AI · All Four Core Modules Have Reached Production-Grade Accuracy

Synthesizability Prediction >90% Average # of Experiments: 1.19

High-Pressure Separation | LCMS Accuracy: 95% (High Confidence: 98%) | NMR >70% Auto-Interpreted

Agentic System · 7 Specialized Agents Have Formed a Complete Digital Chemistry Team

①Project Coordination

②Experiment Setup and Scheduling

③LCMS Reaction Success/Failure Determination

④Purification Work Order Tracking

⑤Quality Control and Structure Confirmation

⑥Delivery Packaging

⑦Material receipt and inventory shortage alerts | Full lifecycle from project initiation to shipping

Platform Achievements

8000/molecules synthesized per month

Reached 5,000 by the end of 2025; increased to 8,000 by April 2026, maintaining stable operations

95%+ / 90%+

Success rate for simple reactions / success rate for complex reactions

Significant Reduction in Total Experiment Volume

More valid data obtained with fewer experiments

Hundreds of Experiments Run in Parallel by a Single Person

A single chemist has shifted from "running reactions one by one" to "batch-managed execution," achieving an order-of-magnitude increase in productivity

AI4S Intelligent Solutions: Agentic HTE (Intelligent Autonomous High-Throughput Experimentation); Building Industry-Leading Autonomous Laboratory Solutions

Reaction Condition Optimization Cycle Greatly Shortened

3-4 Weeks
Traditional HTE



6 Days
Agentic HTE

01 Intent Understanding

Convert experimental objectives and constraints into executable tasks, reducing manual decomposition and repeated back-and-forth communication.

02 Skill Orchestration

Automatically orchestrate condition matrices, robotic steps, and analysis pipelines according to screening objectives.

03 Long-Task Governance

Multi-day experimental tasks are trackable, interruptible, and resumable, ensuring high-throughput workflows run continuously without interruption.

04 Human-Machine Collaboration

Preserve chemist confirmation at key nodes, focusing human effort on judgment rather than repetitive operations.



New Materials & Consumer Health

Advanced Materials: Extending AI4S Core Capabilities to Industrial Cornerstones, Enabling Cross-Scenario Reuse and Cross-Industry Transformation

01 Perspectives & Strategic Focus

- Advanced materials = industry foundation
- Dedicated team + DMTA closed-loop transfer
- Utilizing drug R&D capabilities in advanced materials scenario

02 Tackling PV Challenges Through Collaboration

- Building an automated production line with JinkoSolar
- Advancing advanced materials for perovskite tandem solar cells
- Continuously improving photovoltaic conversion efficiency and cell stability

03 Multiple Order-of-Magnitude Leap in R&D Efficiency

- AI + automated laboratories drive formulation R&D
- Literature mining · Proprietary databases · AI recommendations · Automated validation
- Overall R&D cycle: 4–6 years → 1–6 months

In parallel, XtalPi is expanding into directions such as polymer composites R&D to further broaden the scope of its advanced materials business

New Materials: Intelligent R&D Platform Speed New Materials Discovery, with Major Progress on Perovskite Tandem Cells

AI-Driven Formulation R&D Platform

- The project has built an AI-driven formulation R&D platform centered on literature data mining, dedicated database construction, LLM-based parameter recommendation and automated experimental validation

Performance Validated Devices

- 15mm×15mm module: lab-tested efficiency 27.0%;; third-party certified efficiency 26.51%
- 105mm×105mm module: lab-tested efficiency 23.0%;; third-party certified efficiency 22.74%

Optimization Cycle Compressed from Months to Hours

- The robotic experimental system has completed large-scale sample preparation and built a high-throughput automated tandem cell line, with a designed daily throughput of no less than 1,000 pieces.
- Compared with traditional manual R&D, the single-round optimization cycle has been compressed from months to hours.



Automated Laboratory for Perovskite Cell Preparation & Characterization

Overall R&D Cycle

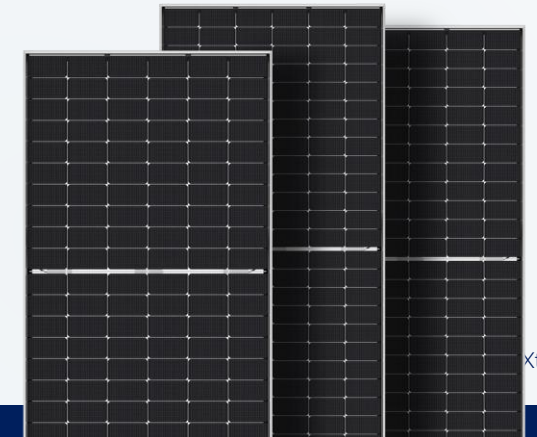
4-6 Years

Traditional R&D Cycle



1-6 Months

Orders-of-Magnitude Leap in R&D Efficiency



Consumer Health: Groland Starts Omni-Channel Build-out in the Large Global Anti-Hair-Loss Market, with Commercialization Moving Ahead

XtalPi Partnership



TOP 1

Tmall Green Bottle Scalp Revitalizing Essence "Anti-Hair Loss, Scalp, Oil New Product Ranking"

TOP 3

During 618, the store ranked among the top 3 new personal-care stores

All-channel sales version

Expected to launch by year-end; two molecules have completed China's new cosmetic-ingredient filing



Groland高岚
头皮精华系列全新上市



Essence
Series

Groland高岚推出铂肽... 华，极致激活毛囊活力，速充头皮胶原，防脱育发，焕活头皮重赋生机。高阶精华适合每日全头皮养护，滚珠精华便于局部使用。另有头皮焕活精华，搭载AI智研双子星科技，多维度实现抗衰、防脱、强韧维稳，精准焕活头皮，为秀发健康注入全新动力。

XtalPi Science Strategic Progression: Evolving into an Open Scientific Intelligence Infrastructure Platform

In July 2026, XtalPi officially launched the XtalPi Science, its scientific intelligence platform, and the Genius Agent matrix of AI agents. As the world's first integrated AI4S platform combining large language models, scientific AI agents, and large-scale robotic experimentation, XtalPi Science standardizes and platformizes an integrated R&D infrastructure that has been validated through real-world projects, lowering the barrier to accessing advanced research capabilities.



Genius Agents

Sign in to your account

USERNAME

PASSWORD

Sign In

or

 Sign in with Feishu

Don't have an account? [Sign up](#)

Intelligent Agents, Accelerated Discovery

Orchestrate autonomous AI agents to streamline your
end to end research



2026.7 XtalPi Science 发布会



5

Financial Results

2026 H1 Financial Key Numbers

RMB mn	2026 H1	2025 H1	YoY
Revenue	394	517	(24%)
Revenue (Excl. Major Pipeline BD)	264	152	+74%
R&D Expenses	368	222	+66%
(Loss) / Profit for the Period	(225)	76	N/A
Adjusted Net (Loss) / Profit	(106)	142	N/A

Cash Balance*: As of June 30, 2026, the Company's total cash balance amounted to RMB 8,671.2 million.

*Cash balance includes our cash and cash equivalents, bank deposits, the current portion of financial assets measured at FVTPL and restricted cash as of 30 June 2026

Excluding Major Pipeline BD: Organic Revenue Grew Substantially, Driving a Significant Narrowing of the Loss Margin

Organic Revenue Grew Substantially

RMB mn

264

152

YoY
+74%¹

2025 H1

2026 H1

Loss Ratio² Narrowed Substantially

RMB mn

2025 H1

Improvement
+58%

2026 H1

(147%)

(89%)

RMB mn

Adjusted net profit / (loss)

Adjusted net profit / (loss) (excluding major pipeline BD)

2025 H1

142

(224)

2026 H1

(106)

(235)

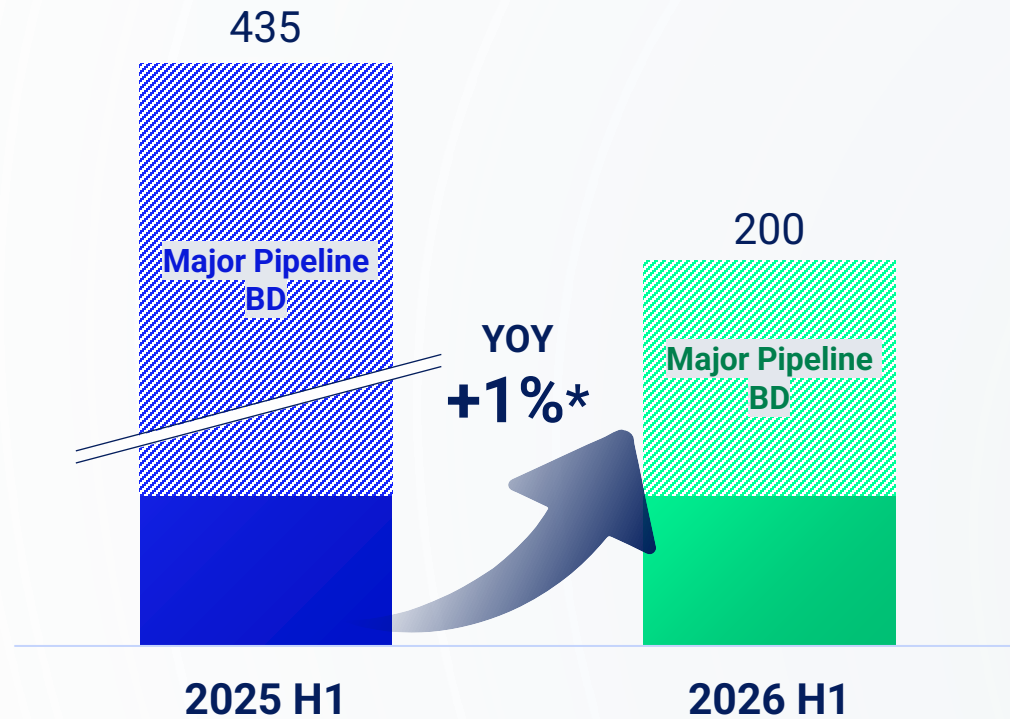
¹Year-on-year revenue growth rate excluding major pipeline BD

²Loss ratio = Non-IFRS net loss excluding major pipeline BD ÷ revenue excluding major pipeline BD

Drug Discovery Solutions Achieves a Strategic Leap, with AI4S Intelligent Solutions Revenue Seeing Exponential Growth

Drug Discovery Solutions

RMB mn



AI4S Intelligent Solutions

RMB mn

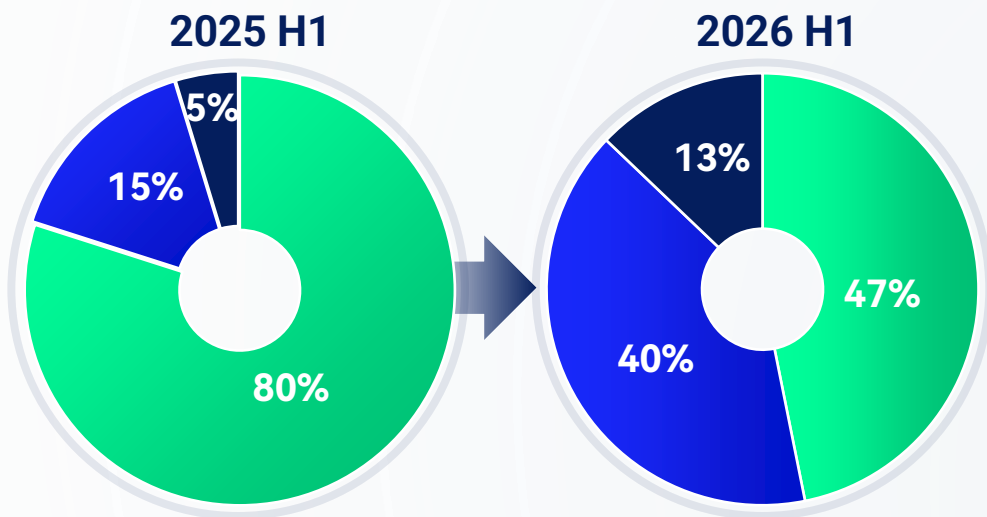


"1%*" refers to the "year-on-year growth rate of revenue from drug discovery solutions excluding major pipeline BD"

Revenue Breakdown by Geography & Business Segment

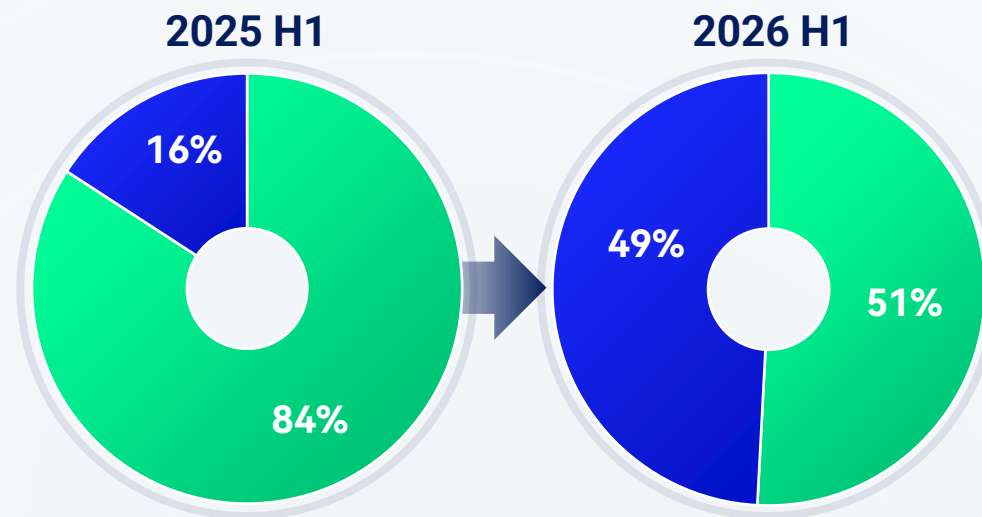
Revenue by Geographical Location (%)

■ United States
 ■ Mainland China
 ■ Others



Revenue by Business Segment (%)

■ Drug Discovery Solutions
 ■ AI4S Intelligent Solutions

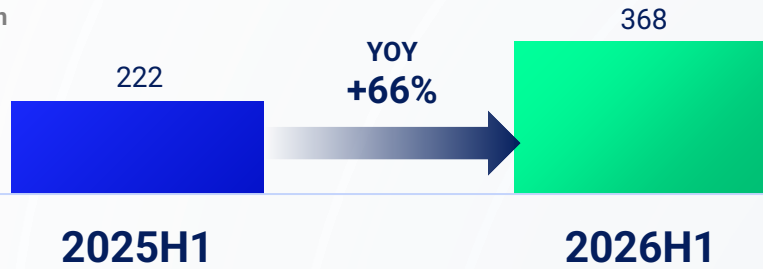


In the first half of 2026, excluding major pipeline BD, revenue across all regions achieved strong year-on-year growth. By business segment, the revenue contribution AI4S Intelligent Solutions increased significantly year on year, with platform value being realized at an accelerated pace and revenue scale achieving leapfrog growth.

Year-on-Year Changes in Operating Expenses

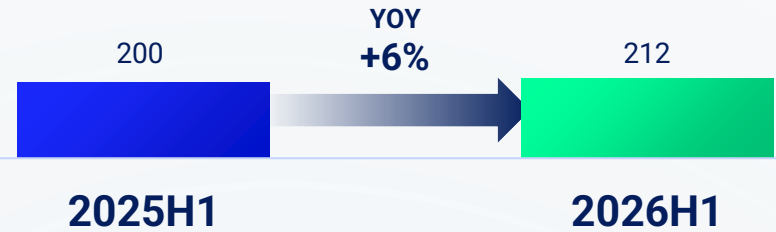
R&D Expenses

RMB mn



G&A Expenses

RMB mn



S&M Expenses

RMB mn



Cost of Revenues

RMB mn



Ample Cash Reserves



The increase in cash balance is primarily attributed to the issuance of convertible bonds, which further strengthened financial reserves.

