

Emotion-Based MCTS Reinforcement Learning System - Detailed Design Specification

Overview

This system is an MCTS-based reinforcement learning algorithm using emotion tokens. While traditional MCTS explores concrete actions a , this approach achieves generalization by exploring abstract emotions e .

1. MCTS Algorithm

1.1 Basic Structure

Each node n maintains the following information:

$$n = \{N(n), W(n), Q(n), P(n), \text{children}(n)\}$$

- $N(n)$: Visit count
- $W(n)$: Cumulative value
- $Q(n) = W(n)/N(n)$: Average value
- $P(n)$: Prior probability (Policy head output)
- $\text{children}(n)$: Set of child nodes

1.2 UCB Selection Formula

At node n , the next emotion e is selected as:

$$e^* = \arg \max_{e \in \mathcal{E}} \left[Q(n, e) + c_{\text{puct}} \cdot P(n, e) \cdot \frac{\sqrt{N(n)}}{1 + N(n, e)} \right]$$

Where:

- $\mathcal{E} = \{0, 1, \dots, 7\}$: Emotion set (8 types)
- $Q(n, e)$: Average value of emotion e
- $P(n, e)$: Prior probability of emotion e
- $N(n, e)$: Visit count of emotion e
- c_{puct} : Exploration coefficient (default: 1.0)

1.3 MCTS Search Flow

Input: Current sequence $s = [s_0, a_0, e_0, r_0, \dots, s_t]$

Output: Visit distribution $\pi = [N(e_0), N(e_1), \dots, N(e_7)]$ (normalized)

Algorithm

```
function MCTS_SEARCH(model, sequence, n_simulations):  
    # 1. Create root node  
    root ← CREATE_NODE()  
  
    # 2. Expand root node  
    outputs ← model.forward(sequence)  
    root.P ← outputs.policy # [8]  
    root.children ← {}  
  
    for e in {0, 1, ..., 7}:  
        child ← CREATE_NODE()  
        child.parent ← root  
        child.emotion ← e  
  
        # Store model predictions  
        child.action_logits ← outputs.actions[e] # [4]  
        child.predicted_state ← outputs.dynamics[e] # [16]  
        child.value ← outputs.values[e] # scalar  
  
        root.children[e] ← child  
  
    # 3. Simulations  
    for i in 1 to n_simulations:  
        SIMULATE(root, sequence, model)  
  
    # 4. Compute visit distribution  
    visit_counts ← [N(root.children[e]) for e in {0, ..., 7}]  
     $\pi$  ← NORMALIZE(visit_counts)  
  
    return  $\pi$ 
```

1.4 Simulation Procedure

```
function SIMULATE(root, real_sequence, model):  
    node ← root  
    path ← [] # Search path  
    imagined_state ← real_sequence.last_state # Imagined starting state  
    depth ← 0  
    MAX_DEPTH ← 3
```

```

# === Selection Phase ===
while NOT node.is_leaf() AND depth < MAX_DEPTH:
    # UCB selection
    e ← SELECT_CHILD_UCB(node)

    path.append((node, e))

    # Get next imagined state
    imagined_state ← node.children[e].predicted_state

    node ← node.children[e]
    depth ← depth + 1

# === Expansion Phase ===
if node.visits > 0 AND depth < MAX_DEPTH:
    # Construct imagined sequence
    imagined_seq ← APPEND_IMAGINED_STATE(real_sequence, imagined_state)

    # Evaluate with model
    outputs ← model.forward(imagined_seq)

    # Expand node
    for e in {0, ..., 7}:
        child ← CREATE_NODE()
        child.parent ← node
        child.action_logits ← outputs.actions[e]
        child.predicted_state ← outputs.dynamics[e]
        child.value ← outputs.values[e]
        node.children[e] ← child

    # Leaf value (average Value across all emotions)
    leaf_value ← MEAN(outputs.values)
else:
    # If unexpanded, use parent's predicted Value
    if path is not empty:
        parent, e ← path[-1]
        leaf_value ← parent.children[e].value
    else:
        leaf_value ← 0.0

# === Backpropagation Phase ===
for (node, e) in REVERSED(path):
    child ← node.children[e]
    child.N ← child.N + 1

```

```
child.W ← child.W + leaf_value  
child.Q ← child.W / child.N
```

1.5 Imagined Sequence Construction

```
function APPEND_IMAGINED_STATE(real_seq, imagined_state):  
    # Copy real sequence  
    new_seq ← COPY(real_seq)  
  
    # Add imagined state (with dummy a, e, r)  
    new_seq.states.append(imagined_state)  
    new_seq.actions.append(0)    # Dummy  
    new_seq.emotions.append(0)  # Dummy  
    new_seq.rewards.append(0.0) # Dummy  
  
    return new_seq
```

1.6 Important Invariants

MCTS imagined sequences are NOT added to the true sequence

```
# During simulation  
imagined_seq = real_seq + [imagined_state, dummy_a, dummy_e, dummy_r] # Temporary  
  
# After simulation, discarded  
  
# Only actual experience is added to true sequence  
e_selected ← ARGMAX(visit_counts)  
a_selected ← model.action_heads[e_selected](state)  
s_next, r_next ← env.step(a_selected)  
  
real_seq ← real_seq + [a_selected, e_selected, r_next, s_next] # Only this persists
```

2. Learning Algorithm

2.1 Four Outputs and Their Roles

The model outputs the following in a single forward pass:

1. **Policy** $\pi_{\theta}(e|s)$: Emotion probability distribution $[8]$
2. **Value** $V_{\theta}^e(s)$: Expected value of emotion e $[8 \times 1]$
3. **Dynamics** $Z_{\theta}^e(s, a^e)$: Next state prediction for emotion e $[8 \times 16]$
4. **Action** $A_{\theta}^e(s)$: Action logits for emotion e $[8 \times 4]$

2.2 Learning Objective Functions

Each output is trained with a different objective function:

2.2.1 Policy Learning (Output 1)

Objective: Imitate MCTS visit distribution

$$\mathcal{L}_\pi = - \sum_{e=0}^7 \pi_{\text{MCTS}}(e) \log \pi_\theta(e|s)$$

Where:

- $\pi_{\text{MCTS}}(e) = N(e) / \sum_{e'} N(e')$: MCTS visit distribution (normalized)
- $\pi_\theta(e|s)$: Model's Policy output (softmax applied)

Gradient: Policy head and Transformer

$$\nabla_{\theta_\pi, \theta_{\text{transformer}}} \mathcal{L}_\pi$$

2.2.2 Value Learning (Output 2)

Objective: Predict actual returns

$$\mathcal{L}_V = (V_\theta^{e_{\text{selected}}}(s) - G_{\text{actual}})^2$$

Where:

- e_{selected} : Emotion selected by MCTS
- $G_{\text{actual}} = \sum_{t'=t}^T \gamma^{t'-t} r_{t'}$: Actual discounted return
- $\gamma = 0.99$: Discount factor

Gradient: Value head and Transformer

$$\nabla_{\theta_V, \theta_{\text{transformer}}} \mathcal{L}_V$$

2.2.3 Dynamics Learning (Output 3)

Objective: Predict next actual state

$$\mathcal{L}_Z = \|Z_\theta^{e_{\text{selected}}}(s, a^{e_{\text{selected}}}) - s_{\text{next}}^{\text{actual}}\|_2^2$$

Where:

- a^{selected} : Action head output (**frozen, no gradient**)
- $s_{\text{next}}^{\text{actual}}$: Actual next state from environment

Important: Action head is frozen

$$a^{\text{selected}} = \text{stopgrad}(A_{\theta}^{\text{selected}}(s))$$

Gradient: Dynamics head and Transformer

$$\nabla_{\theta_Z, \theta_{\text{transformer}}} \mathcal{L}_Z$$

2.2.4 Action Learning (Output 4)

Objective: Reproduce the state predicted by Dynamics

$$\mathcal{L}_A = \|Z_{\theta}^{\text{selected}}(s, a^{\text{selected}}) - z_{\text{target}}\|_2^2$$

Where:

- $a^{\text{selected}} = A_{\theta}^{\text{selected}}(s)$: Action head output (**trainable**)
- $z_{\text{target}} = \text{stopgrad}(Z_{\theta}^{\text{selected}}(s, a^{\text{selected}}))$: Dynamics prediction (**frozen**)

Important: Dynamics head is frozen, only Action is trained

Gradient: Action head and Transformer

$$\nabla_{\theta_A, \theta_{\text{transformer}}} \mathcal{L}_A$$

2.3 Alternating Updates

The four losses are optimized **alternately**:

for each timestep t in episode:

$\text{sequence_t} \leftarrow [s_0, a_0, e_0, r_0, \dots, s_t]$

$e_{\text{selected}} \leftarrow \text{emotions}[t]$

=== Phase 1: Dynamics update ===

$\text{outputs} \leftarrow \text{model.forward}(\text{sequence_t})$

$a_e \leftarrow \text{stopgrad}(\text{outputs.actions}[e_{\text{selected}}])$ # Action frozen

$h_t \leftarrow \text{outputs.hidden}$

$z_{\text{pred}} \leftarrow \text{model.dynamics_heads}[e_{\text{selected}}](h_t, a_e)$

$s_{\text{actual}} \leftarrow \text{states}[t+1]$

$L_{\text{dynamics}} \leftarrow \text{MSE}(z_{\text{pred}}, s_{\text{actual}})$

$\text{optimizer_dynamics.zero_grad}()$

$L_{\text{dynamics.backward}}()$

$\text{optimizer_dynamics.step}()$

=== Phase 2: Action update ===

$\text{outputs} \leftarrow \text{model.forward}(\text{sequence_t})$

$a_e \leftarrow \text{outputs.actions}[e_{\text{selected}}]$ # Action trainable

$h_t \leftarrow \text{stopgrad}(\text{outputs.hidden})$

$z_{\text{target}} \leftarrow \text{stopgrad}(\text{outputs.dynamics}[e_{\text{selected}}])$ # Dynamics frozen

$z_{\text{pred}} \leftarrow \text{model.dynamics_heads}[e_{\text{selected}}](h_t, a_e)$

$L_{\text{action}} \leftarrow \text{MSE}(z_{\text{pred}}, z_{\text{target}})$

$\text{optimizer_action.zero_grad}()$

$L_{\text{action.backward}}()$

$\text{optimizer_action.step}()$

=== Phase 3: Value update ===

$\text{outputs} \leftarrow \text{model.forward}(\text{sequence_t})$

$v_{\text{pred}} \leftarrow \text{outputs.values}[e_{\text{selected}}]$

$G_{\text{actual}} \leftarrow \text{returns}[t]$

$L_{\text{value}} \leftarrow \text{MSE}(v_{\text{pred}}, G_{\text{actual}})$

$\text{optimizer_value.zero_grad}()$

$L_{\text{value.backward}}()$

$\text{optimizer_value.step}()$

```
# === Phase 4: Policy update ===
outputs ← model.forward(sequence_t)

π_pred ← outputs.policy
π_mcts ← mcts_distributions[t]

L_policy ← -SUM(π_mcts * LOG(π_pred + ε))

optimizer_policy.zero_grad()
L_policy.backward()
optimizer_policy.step()
```

2.4 Gradient Flow

Phase 1: Dynamics Learning

```
s_actual → MSE → Dynamics Head → Transformer → backward
          ↑
        Action Head (frozen, no grad)
```

Phase 2: Action Learning

```
z_target (frozen) → MSE → Dynamics Head (frozen)
                  ↑
        Action Head → Transformer → backward
```

Phase 3: Value Learning

```
G_actual → MSE → Value Head → Transformer → backward
```

Phase 4: Policy Learning

```
π_mcts → CrossEntropy → Policy Head → Transformer → backward
```

Important: Gradients flow to the Transformer in all Phases. This optimizes the history encoding h_t for each task.

2.5 Important Design Decisions

Why Dynamics Takes Action as Input

$$Z_{\theta}(s, a) = f_{\theta}(\text{concat}(h(s), a))$$

Reasons:

1. Action gradients flow through Dynamics (Phase 2)
2. Dynamics is conditioned on Action ($p(s_{t+1}|s_t, a_t)$)
3. Dynamics can be used as a "differentiable environment model" during Action learning

Incorrect Implementation:

```
python

# ✗ This doesn't propagate gradients
z_pred = dynamics_head(h_t) # Doesn't take a
a = action_head(h_t)
loss = MSE(z_pred, target) # No gradient to a
```

Correct Implementation:

```
python

# ✓ Gradients flow
a = action_head(h_t)
z_pred = dynamics_head(h_t, a) # Takes a as input
loss = MSE(z_pred, target) # Gradients flow to a
```

Why Alternating Updates

Reasons:

1. **Phase 1:** Dynamics learns reality → Accurate world model
2. **Phase 2:** Action utilizes Dynamics → Better action selection
3. Mutual improvement feedback loop
4. **Transformer learns from all four tasks** → Multi-task learning effect

Regular end-to-end learning can be unstable when Dynamics and Action are updated simultaneously.

About Four Transformer Updates:

At each timestep, the Transformer is updated four times (once per Phase). This has the following advantages:

1. Learns from multiple objective functions → More general representation
2. Doesn't overspecialize to each task
3. Learns h_t useful for Policy/Value/Dynamics/Action

However, there's a risk of overfitting without proper learning rate settings. Recommended:

- Head learning rate: 3×10^{-4}

- Transformer learning rate: 1×10^{-4} (lower than Heads)

Alternatively, you can share Transformer parameters across all Optimizers, effectively making it a single update (implementation-dependent).

3. Important Implementation Points

3.1 Sequence Management

```
python

# ✓ Correct: Store only actual experience
real_sequence = [s_0, a_0, e_0, r_0, s_1, a_1, e_1, r_1, ...]

# ✗ Wrong: Store MCTS imagination
# real_sequence += imagined_transitions # Don't do this
```

3.2 Optimizer Configuration

Important: Transformer is shared across all Optimizers

```
python

# ✓ Correct: Optimize each Head with Transformer
optimizer_dynamics = Adam(
    list(dynamics_heads.parameters()) + list(transformer.parameters())
)
optimizer_action = Adam(
    list(action_heads.parameters()) + list(transformer.parameters())
)
optimizer_value = Adam(
    list(value_heads.parameters()) + list(transformer.parameters())
)
optimizer_policy = Adam(
    list(policy_head.parameters()) + list(transformer.parameters())
)

# ✗ Wrong 1: Transformer not trained
# optimizer_dynamics = Adam(dynamics_heads.parameters()) # No Transformer

# ✗ Wrong 2: Single Optimizer
# optimizer = Adam(model.parameters()) # Can't do Alternating Updates
```

Note: The Transformer is updated four times (once per Phase). This is intentional design, acquiring general representations by learning from multiple tasks.

3.3 Saving MCTS Visit Distribution

```
python

# ✓ Correct: Save during episode collection
episode = {
    'states': [...],
    'actions': [...],
    'emotions': [...],
    'rewards': [...],
    'mcts_distributions': [...] # This is essential
}

# ✗ Wrong: Don't save visit distribution
# → Can't train Policy
```

3.4 Temperature Scheduling for Emotion Selection

```
python

# Early training: High temperature (exploration-focused)
temperature = 1.0

# Late training: Low temperature (exploitation-focused)
temperature = 0.1

# Sampling
if temperature == 0:
    e_selected = argmax(visit_dist)
else:
    probs = visit_dist ** (1 / temperature)
    probs = probs / sum(probs)
    e_selected = random.choice(probs)
```

4. Mathematical Details

4.1 Derivation of UCB Formula

The UCB formula combines two terms:

Exploitation term:

$$Q(n, e) = \frac{W(n, e)}{N(n, e)}$$

Exploration term:

$$U(n, e) = c_{\text{puct}} \cdot P(n, e) \cdot \frac{\sqrt{N(n)}}{1 + N(n, e)}$$

This maximizes the upper bound:

$$\text{UCB}(n, e) = Q(n, e) + U(n, e)$$

As $N(n, e) \rightarrow \infty$, $U(n, e) \rightarrow 0$, and exploitation dominates.

4.2 Justification for Policy Loss

The MCTS visit distribution π_{MCTS} approximates the optimal policy with sufficient simulations. The Policy network distills this distribution:

$$\min_{\theta} D_{\text{KL}}(\pi_{\text{MCTS}} \parallel \pi_{\theta}) = - \sum_e \pi_{\text{MCTS}}(e) \log \pi_{\theta}(e) + \text{const}$$

This enables the Policy network to "select good emotions without MCTS".

4.3 Theory of Dynamics-Action Separated Learning

Goal: Modeling $p(s_{t+1} | s_t, a_t)$

Dynamics learns the laws of the world:

$$Z_{\theta}(s_t, a_t) \approx s_{t+1}$$

Action is optimized using Dynamics as the environment:

$$\max_a V(Z_{\theta}(s_t, a))$$

This is a form of model-based RL; if Dynamics is accurate, Action also improves.

5. Overall Algorithm Flow

```
# === Training Loop ===  
for episode in range(num_episodes):  
    env.reset()  
    sequence = [initial_state]  
    episode_data = {states: [], actions: [], emotions: [], rewards: [], mcts_dists: []}
```

while not done:

1. Run MCTS

visit_dist = MCTS_SEARCH(model, sequence, n_simulations=50)

2. Select emotion

e = SAMPLE(visit_dist, temperature)

3. Get action

outputs = model.forward(sequence)

a_logits = outputs.actions[e]

a = SAMPLE(softmax(a_logits))

4. Execute in environment

s_next, r, done = env.step(a)

5. Store experience

episode_data.states.append(current_state)

episode_data.actions.append(a)

episode_data.emotions.append(e)

episode_data.rewards.append(r)

episode_data.mcts_dists.append(visit_dist)

6. Update sequence (only actual experience)

sequence += [a, e, r, s_next]

current_state = s_next

7. Compute returns

episode_data.returns = COMPUTE_RETURNS(episode_data.rewards, gamma=0.99)

8. Alternating Updates

for t in range(len(episode_data.states) - 1):

 TRAIN_STEP(model, episode_data, t, optimizers)

9. Save to buffer

buffer.add(episode_data)

6. Checklist

Items to verify during implementation:

- ☐ MCTS selects nodes using UCB formula
- ☐ MCTS returns visit distribution
- ☐ MCTS imagined sequences are not added to true sequence
- ☐ Dynamics Head takes (h_t, a) as input

- ☐ Phase 1: stopgrad Action
 - ☐ Phase 2: stopgrad Dynamics, Action is trainable
 - ☐ Phase 3: Train Value with actual returns
 - ☐ Phase 4: Train Policy with MCTS visit distribution
 - ☐ Four independent Optimizers
 - ☐ Save MCTS visit distribution to episode data
 - ☐ Implement temperature scheduling
-

This is the implementation-level detailed design specification.